

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
 Data File : BF095726.D  
 Acq On : 7 Jun 2017 00:16  
 Operator : SJ/MA  
 Sample : I3482-07  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CL-01-060517-C

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.587       | 506           | 510         | 529          | rBV2     | 46688          | 126800        | 7.73%           | 0.518%        |
| 2         | 5.269       | 622           | 626         | 651          | rBV      | 842606         | 1449482       | 88.33%          | 5.920%        |
| 3         | 6.940       | 905           | 910         | 922          | rBV      | 939124         | 1406274       | 85.70%          | 5.743%        |
| 4         | 7.034       | 922           | 926         | 938          | rVB      | 1133698        | 1633471       | 99.55%          | 6.671%        |
| 5         | 7.381       | 981           | 985         | 1000         | rVB      | 343603         | 429972        | 26.20%          | 1.756%        |
| 6         | 7.616       | 1020          | 1025        | 1038         | rBV      | 909261         | 1196893       | 72.94%          | 4.888%        |
| 7         | 8.298       | 1136          | 1141        | 1159         | rBV      | 626859         | 873527        | 53.23%          | 3.568%        |
| 8         | 9.410       | 1326          | 1330        | 1343         | rBV      | 446491         | 581130        | 35.41%          | 2.373%        |
| 9         | 10.586      | 1526          | 1530        | 1539         | rVB      | 21707          | 30847         | 1.88%           | 0.126%        |
| 10        | 10.733      | 1550          | 1555        | 1563         | rBV      | 21137          | 31545         | 1.92%           | 0.129%        |
| 11        | 11.192      | 1628          | 1633        | 1647         | rVB      | 1389368        | 1640916       | 100.00%         | 6.702%        |
| 12        | 11.716      | 1717          | 1722        | 1726         | rBV      | 38241          | 52438         | 3.20%           | 0.214%        |
| 13        | 11.757      | 1726          | 1729        | 1738         | rVB3     | 25041          | 37750         | 2.30%           | 0.154%        |
| 14        | 11.886      | 1746          | 1751        | 1759         | rBV      | 156005         | 199723        | 12.17%          | 0.816%        |
| 15        | 12.010      | 1768          | 1772        | 1780         | rVB      | 64113          | 96371         | 5.87%           | 0.394%        |
| 16        | 12.233      | 1802          | 1810        | 1815         | rBV2     | 453630         | 555081        | 33.83%          | 2.267%        |
| 17        | 12.280      | 1815          | 1818        | 1828         | rVB3     | 48893          | 73438         | 4.48%           | 0.300%        |
| 18        | 12.604      | 1867          | 1873        | 1878         | rVB2     | 24249          | 45346         | 2.76%           | 0.185%        |
| 19        | 12.669      | 1878          | 1884        | 1888         | rBV2     | 34884          | 47099         | 2.87%           | 0.192%        |
| 20        | 12.780      | 1894          | 1903        | 1909         | rBV3     | 37710          | 74480         | 4.54%           | 0.304%        |
| 21        | 12.927      | 1920          | 1928        | 1930         | rBV3     | 16105          | 29218         | 1.78%           | 0.119%        |
| 22        | 13.092      | 1947          | 1956        | 1959         | rBV3     | 52359          | 90719         | 5.53%           | 0.371%        |
| 23        | 13.127      | 1959          | 1962        | 1972         | rVB      | 51470          | 87817         | 5.35%           | 0.359%        |
| 24        | 13.216      | 1972          | 1977        | 1979         | rBV3     | 19219          | 27960         | 1.70%           | 0.114%        |
| 25        | 13.421      | 2005          | 2012        | 2020         | rBV6     | 22271          | 67175         | 4.09%           | 0.274%        |
| 26        | 13.527      | 2025          | 2030        | 2039         | rVV      | 561720         | 728204        | 44.38%          | 2.974%        |
| 27        | 13.757      | 2066          | 2069        | 2076         | rVB3     | 22752          | 31069         | 1.89%           | 0.127%        |
| 28        | 13.869      | 2083          | 2088        | 2098         | rBV3     | 60647          | 122367        | 7.46%           | 0.500%        |
| 29        | 14.027      | 2109          | 2115        | 2119         | rBV4     | 33111          | 65970         | 4.02%           | 0.269%        |
| 30        | 14.069      | 2119          | 2122        | 2127         | rVB2     | 32446          | 44436         | 2.71%           | 0.181%        |
| 31        | 14.151      | 2127          | 2136        | 2139         | rBV5     | 21078          | 39041         | 2.38%           | 0.159%        |
| 32        | 14.257      | 2149          | 2154        | 2158         | rBV3     | 20317          | 30445         | 1.86%           | 0.124%        |
| 33        | 14.463      | 2186          | 2189        | 2194         | rVB      | 27186          | 35272         | 2.15%           | 0.144%        |
| 34        | 14.598      | 2208          | 2212        | 2213         | rBV      | 32967          | 34777         | 2.12%           | 0.142%        |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
 Data File : BF095726.D  
 Acq On : 7 Jun 2017 00:16  
 Operator : SJ/MA  
 Sample : I3482-07  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CL-01-060517-C

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 14.627 | 2213 | 2217 | 2220 | rVV  | 365758 | 461653  | 28.13% | 1.885% |
| 36 | 14.663 | 2220 | 2223 | 2229 | rVB  | 305539 | 367090  | 22.37% | 1.499% |
| 37 | 14.751 | 2234 | 2238 | 2244 | rVB  | 110765 | 145420  | 8.86%  | 0.594% |
| 38 | 14.857 | 2249 | 2256 | 2259 | rBV2 | 36882  | 57669   | 3.51%  | 0.236% |
| 39 | 15.039 | 2281 | 2287 | 2289 | rBV4 | 16480  | 28522   | 1.74%  | 0.116% |
| 40 | 15.127 | 2300 | 2302 | 2305 | rVV2 | 24580  | 26030   | 1.59%  | 0.106% |
| 41 | 15.163 | 2305 | 2308 | 2313 | rVB  | 56691  | 62728   | 3.82%  | 0.256% |
| 42 | 15.280 | 2323 | 2328 | 2331 | rBV3 | 32883  | 57574   | 3.51%  | 0.235% |
| 43 | 15.427 | 2349 | 2353 | 2357 | rBV  | 168539 | 189918  | 11.57% | 0.776% |
| 44 | 15.474 | 2357 | 2361 | 2367 | rVB  | 182514 | 222308  | 13.55% | 0.908% |
| 45 | 15.545 | 2370 | 2373 | 2376 | rBV  | 86231  | 90626   | 5.52%  | 0.370% |
| 46 | 15.586 | 2376 | 2380 | 2383 | rVV2 | 199064 | 291281  | 17.75% | 1.190% |
| 47 | 15.621 | 2383 | 2386 | 2399 | rVB  | 153970 | 284304  | 17.33% | 1.161% |
| 48 | 15.780 | 2410 | 2413 | 2417 | rBV3 | 38525  | 46729   | 2.85%  | 0.191% |
| 49 | 15.880 | 2426 | 2430 | 2435 | rVV3 | 100353 | 163593  | 9.97%  | 0.668% |
| 50 | 16.051 | 2456 | 2459 | 2463 | rVB3 | 30875  | 35546   | 2.17%  | 0.145% |
| 51 | 16.098 | 2463 | 2467 | 2469 | rBV2 | 54014  | 69111   | 4.21%  | 0.282% |
| 52 | 16.139 | 2471 | 2474 | 2478 | rVV  | 80043  | 92247   | 5.62%  | 0.377% |
| 53 | 16.174 | 2478 | 2480 | 2486 | rVB  | 49638  | 50891   | 3.10%  | 0.208% |
| 54 | 16.239 | 2487 | 2491 | 2495 | rBV  | 229285 | 297079  | 18.10% | 1.213% |
| 55 | 16.280 | 2495 | 2498 | 2502 | rVV4 | 129652 | 183391  | 11.18% | 0.749% |
| 56 | 16.321 | 2502 | 2505 | 2508 | rVB2 | 64712  | 70017   | 4.27%  | 0.286% |
| 57 | 16.362 | 2508 | 2512 | 2517 | rVB  | 138508 | 174848  | 10.66% | 0.714% |
| 58 | 16.445 | 2519 | 2526 | 2531 | rBV  | 519834 | 696406  | 42.44% | 2.844% |
| 59 | 16.492 | 2531 | 2534 | 2538 | rVB2 | 57673  | 72006   | 4.39%  | 0.294% |
| 60 | 16.545 | 2538 | 2543 | 2546 | rBV2 | 91254  | 111198  | 6.78%  | 0.454% |
| 61 | 16.580 | 2546 | 2549 | 2552 | rBV  | 57192  | 59620   | 3.63%  | 0.243% |
| 62 | 16.639 | 2556 | 2559 | 2563 | rBV2 | 33473  | 57154   | 3.48%  | 0.233% |
| 63 | 16.686 | 2563 | 2567 | 2571 | rBV3 | 34051  | 41856   | 2.55%  | 0.171% |
| 64 | 16.745 | 2572 | 2577 | 2583 | rVB  | 793267 | 906164  | 55.22% | 3.701% |
| 65 | 16.821 | 2588 | 2590 | 2595 | rVB3 | 58056  | 78783   | 4.80%  | 0.322% |
| 66 | 16.874 | 2595 | 2599 | 2603 | rBV  | 76174  | 90060   | 5.49%  | 0.368% |
| 67 | 16.939 | 2607 | 2610 | 2615 | rBV4 | 83283  | 101504  | 6.19%  | 0.415% |
| 68 | 17.004 | 2615 | 2621 | 2628 | rVB  | 948617 | 1164469 | 70.96% | 4.756% |
| 69 | 17.074 | 2628 | 2633 | 2636 | rBV2 | 104988 | 165712  | 10.10% | 0.677% |
| 70 | 17.186 | 2647 | 2652 | 2654 | rBV3 | 89139  | 164752  | 10.04% | 0.673% |
| 71 | 17.215 | 2654 | 2657 | 2663 | rVB2 | 166027 | 229215  | 13.97% | 0.936% |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
 Data File : BF095726.D  
 Acq On : 7 Jun 2017 00:16  
 Operator : SJ/MA  
 Sample : I3482-07  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CL-01-060517-C

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

Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |        |        |        |        |
|-----|--------|------|------|------|------|--------|--------|--------|--------|
| 72  | 17.304 | 2663 | 2672 | 2676 | rBV2 | 116201 | 210093 | 12.80% | 0.858% |
| 73  | 17.345 | 2676 | 2679 | 2687 | rBV2 | 183100 | 263349 | 16.05% | 1.076% |
| 74  | 17.409 | 2688 | 2690 | 2694 | rVV2 | 40994  | 43183  | 2.63%  | 0.176% |
| 75  | 17.462 | 2695 | 2699 | 2703 | rVV  | 166293 | 179176 | 10.92% | 0.732% |
| 76  | 17.504 | 2703 | 2706 | 2709 | rVV  | 126495 | 128444 | 7.83%  | 0.525% |
| 77  | 17.645 | 2727 | 2730 | 2735 | rBV5 | 26738  | 42685  | 2.60%  | 0.174% |
| 78  | 17.721 | 2739 | 2743 | 2744 | rBV3 | 35118  | 47107  | 2.87%  | 0.192% |
| 79  | 17.857 | 2763 | 2766 | 2771 | rBV2 | 90443  | 136561 | 8.32%  | 0.558% |
| 80  | 17.898 | 2771 | 2773 | 2777 | rVB  | 68040  | 78764  | 4.80%  | 0.322% |
| 81  | 17.962 | 2782 | 2784 | 2787 | rBV  | 34004  | 37425  | 2.28%  | 0.153% |
| 82  | 18.015 | 2790 | 2793 | 2796 | rVV2 | 86443  | 114699 | 6.99%  | 0.468% |
| 83  | 18.045 | 2796 | 2798 | 2801 | rVV2 | 113024 | 116646 | 7.11%  | 0.476% |
| 84  | 18.080 | 2801 | 2804 | 2807 | rVV  | 30918  | 30191  | 1.84%  | 0.123% |
| 85  | 18.280 | 2829 | 2838 | 2843 | rBV5 | 86250  | 215746 | 13.15% | 0.881% |
| 86  | 18.339 | 2843 | 2848 | 2852 | rVV2 | 550277 | 900642 | 54.89% | 3.678% |
| 87  | 18.374 | 2852 | 2854 | 2861 | rVB  | 336944 | 419010 | 25.54% | 1.711% |
| 88  | 18.468 | 2867 | 2870 | 2874 | rBV2 | 66430  | 76577  | 4.67%  | 0.313% |
| 89  | 18.674 | 2902 | 2905 | 2910 | rVB3 | 88264  | 124340 | 7.58%  | 0.508% |
| 90  | 18.803 | 2924 | 2927 | 2929 | rBV2 | 46176  | 52321  | 3.19%  | 0.214% |
| 91  | 18.851 | 2929 | 2935 | 2939 | rVV  | 157725 | 240310 | 14.64% | 0.981% |
| 92  | 18.892 | 2939 | 2942 | 2945 | rVB  | 106479 | 108154 | 6.59%  | 0.442% |
| 93  | 19.003 | 2958 | 2961 | 2968 | rBV3 | 57024  | 117283 | 7.15%  | 0.479% |
| 94  | 19.062 | 2969 | 2971 | 2975 | rVV2 | 49083  | 53243  | 3.24%  | 0.217% |
| 95  | 19.609 | 3060 | 3064 | 3067 | rBV  | 253197 | 346486 | 21.12% | 1.415% |
| 96  | 19.898 | 3110 | 3113 | 3118 | rBV  | 169733 | 193729 | 11.81% | 0.791% |
| 97  | 19.956 | 3120 | 3123 | 3127 | rVB  | 232618 | 241387 | 14.71% | 0.986% |
| 98  | 20.015 | 3129 | 3133 | 3136 | rBV  | 235207 | 273411 | 16.66% | 1.117% |
| 99  | 21.192 | 3331 | 3333 | 3339 | rVB  | 92412  | 125475 | 7.65%  | 0.512% |
| 100 | 21.527 | 3387 | 3390 | 3398 | rVB  | 87662  | 144026 | 8.78%  | 0.588% |

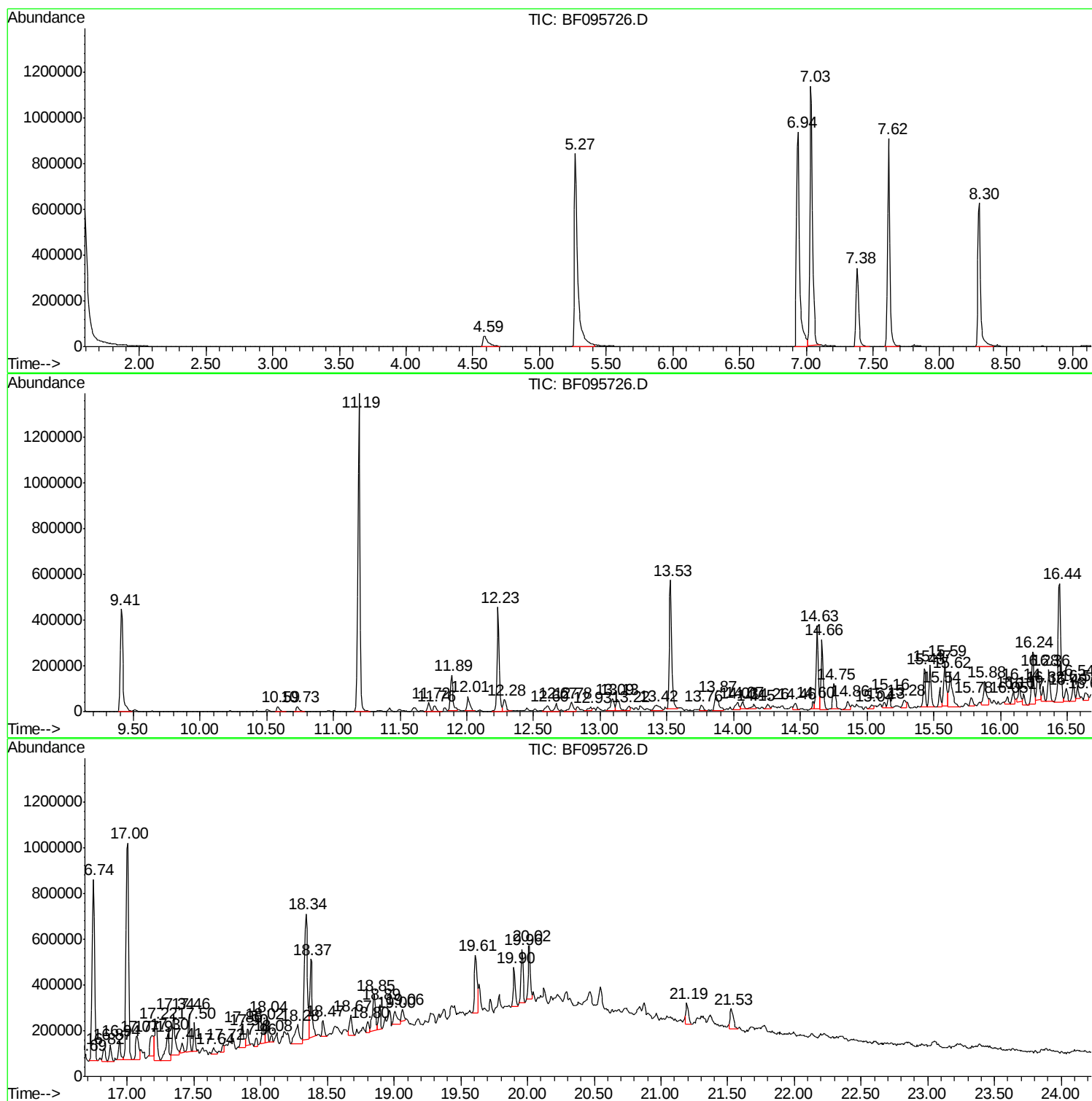
Sum of corrected areas: 24484990

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

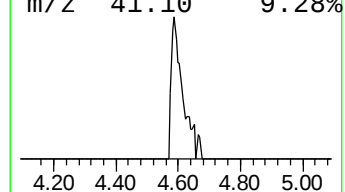
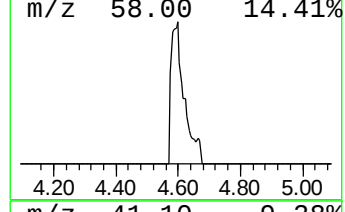
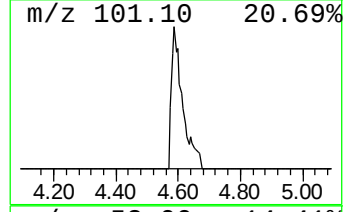
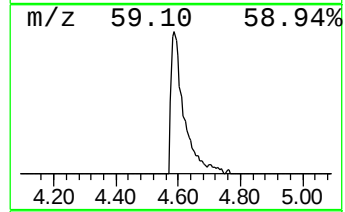
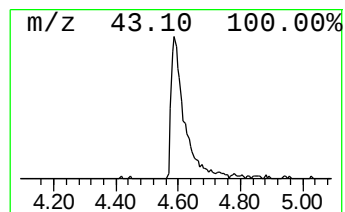
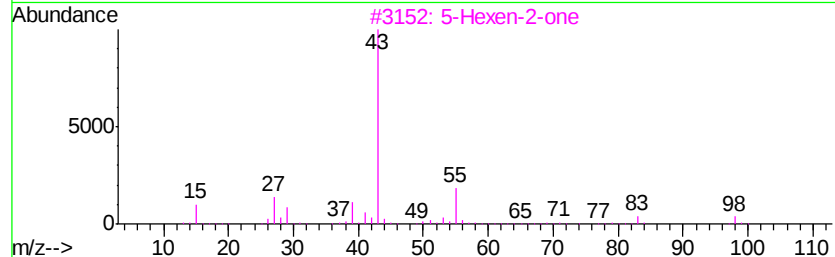
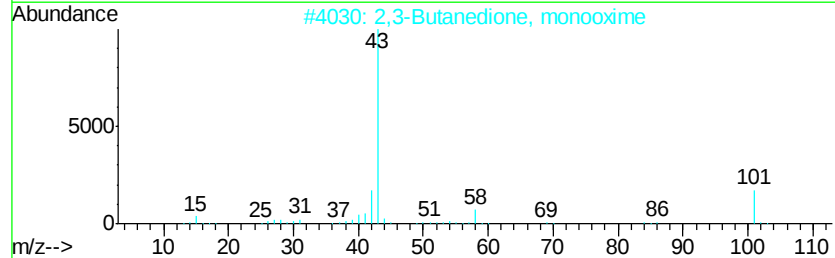
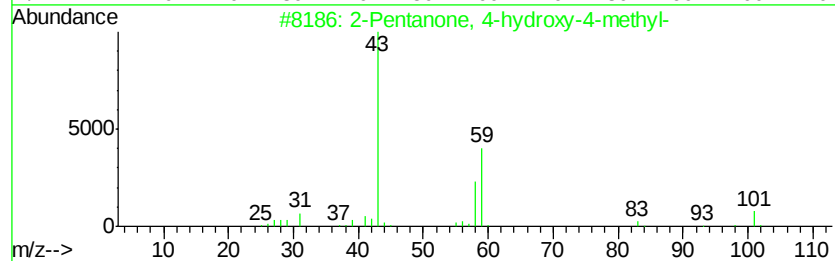
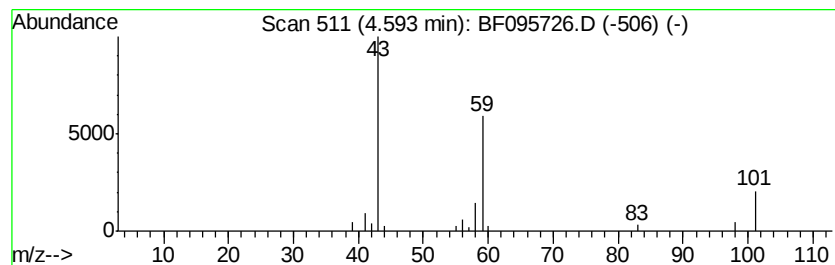
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.59 | 5.90 ng | 126800 | 1,4-Dichlorobenzene-d4 | 7.38 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2 | 56   |
| 2    |    |   | 2,3-Butanedione, monooxime          | 101 | C4H7NO2 | 000057-71-6 | 9    |
| 3    |    |   | 5-Hexen-2-one                       | 98  | C6H10O  | 000109-49-9 | 9    |
| 4    |    |   | 3-Pentanol, 2-methyl-               | 102 | C6H14O  | 000565-67-3 | 9    |
| 5    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

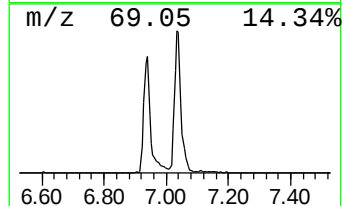
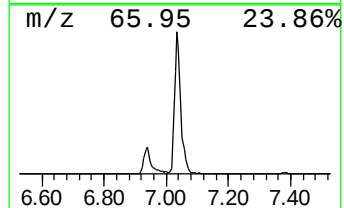
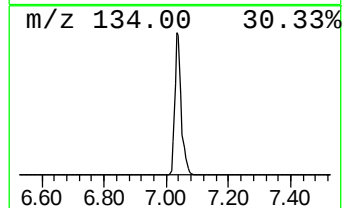
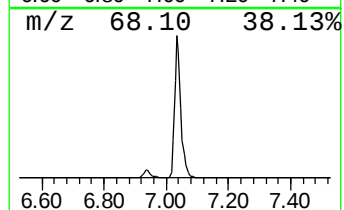
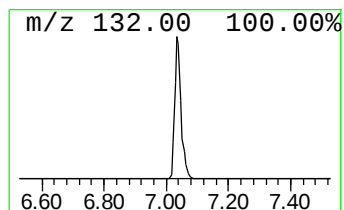
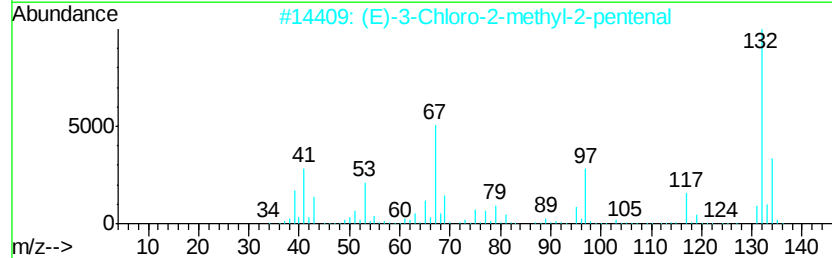
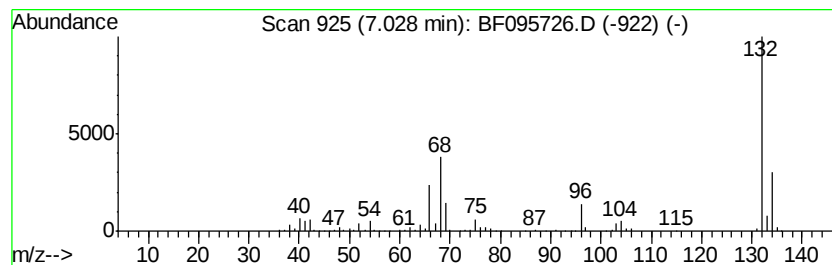
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown7.03 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.03 | 75.98 ng | 1633470 | 1,4-Dichlorobenzene-d4 | 7.38 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 25   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 10   |
| 5    |    | 4-Chloro-2-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-00-5  | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

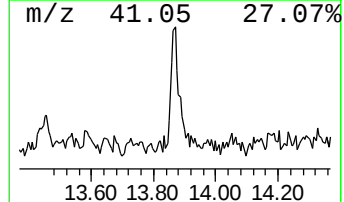
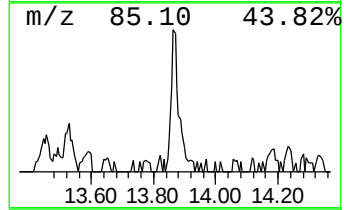
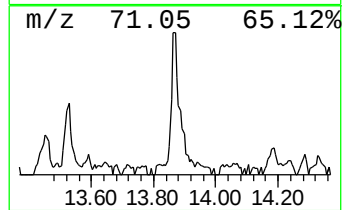
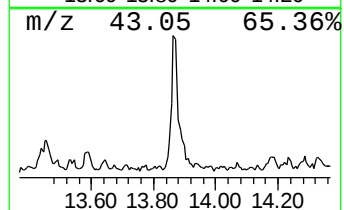
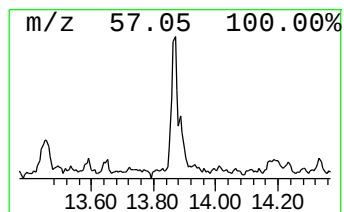
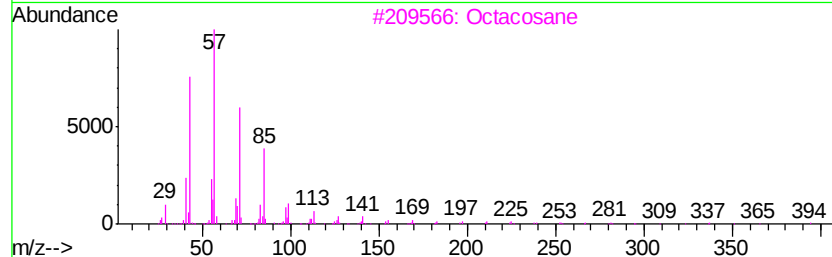
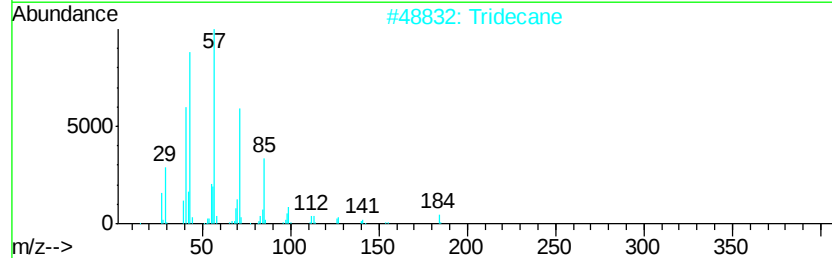
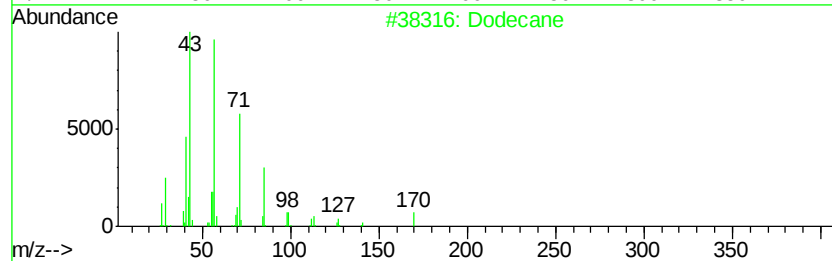
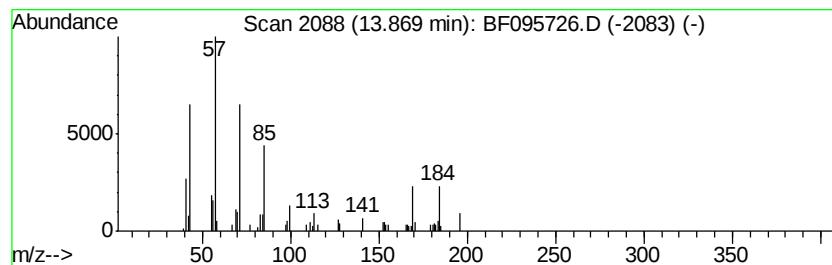
Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Dodecane Concentration Rank 15

| R.T.    | EstConc     | Area         | Relative to ISTD |         | R.T.        |      |
|---------|-------------|--------------|------------------|---------|-------------|------|
| 13.87   | 5.30 ng     | 122367       | Phenanthrene-d10 |         | 14.63       |      |
| Hit# of | 5           | Tentative ID | MW               | MolForm | CAS#        | Qual |
| 1       | Dodecane    |              | 170              | C12H26  | 000112-40-3 | 70   |
| 2       | Tridecane   |              | 184              | C13H28  | 000629-50-5 | 64   |
| 3       | Octacosane  |              | 394              | C28H58  | 000630-02-4 | 62   |
| 4       | Hexadecane  |              | 226              | C16H34  | 000544-76-3 | 58   |
| 5       | Heneicosane |              | 296              | C21H44  | 000629-94-7 | 58   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

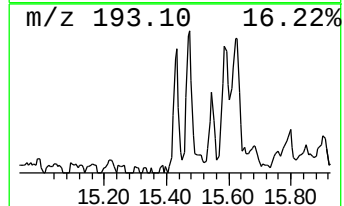
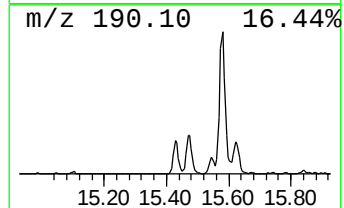
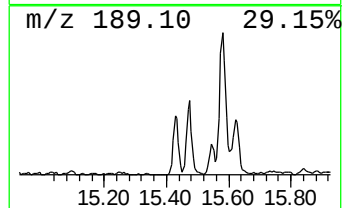
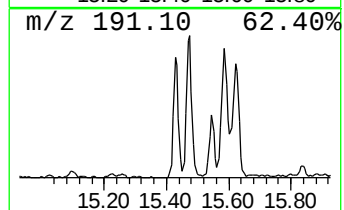
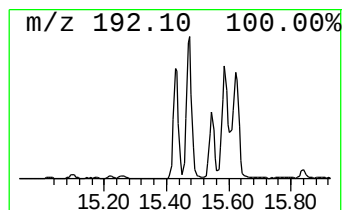
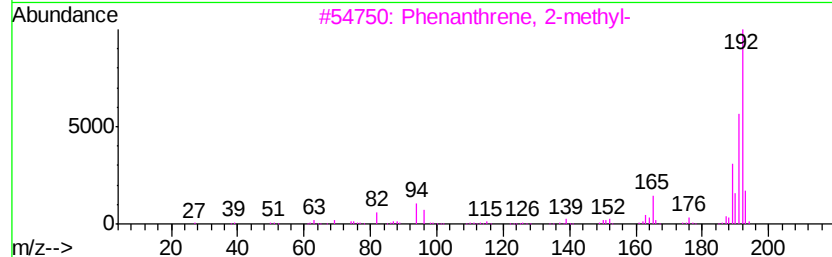
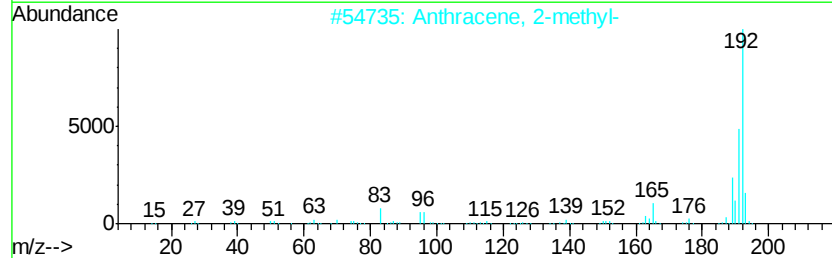
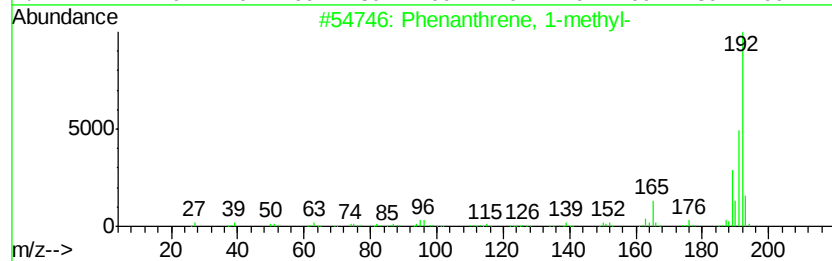
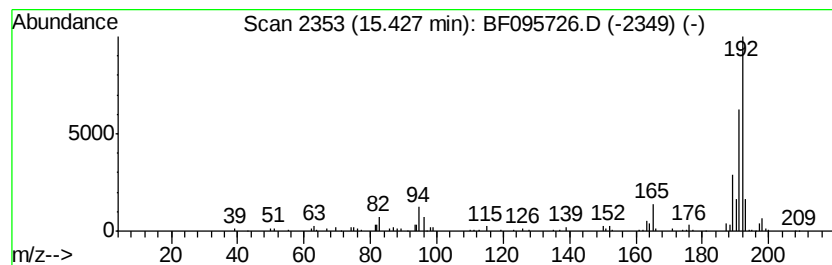
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.43 | 8.23 ng | 189918 | Phenanthrene-d10 | 14.63 |

| Hit# of | 5                                     | Tentative ID | MW     | MolForm     | CAS# | Qual |
|---------|---------------------------------------|--------------|--------|-------------|------|------|
| 1       | Phenanthrene, 1-methyl-               | 192          | C15H12 | 000832-69-9 | 96   |      |
| 2       | Anthracene, 2-methyl-                 | 192          | C15H12 | 000613-12-7 | 93   |      |
| 3       | Phenanthrene, 2-methyl-               | 192          | C15H12 | 002531-84-2 | 91   |      |
| 4       | Anthracene, 1-methyl-                 | 192          | C15H12 | 000610-48-0 | 90   |      |
| 5       | 1H-Cyclopropa[1]phenanthrene, 1a, ... | 192          | C15H12 | 000949-41-7 | 89   |      |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

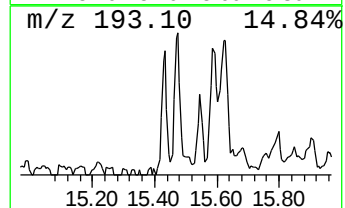
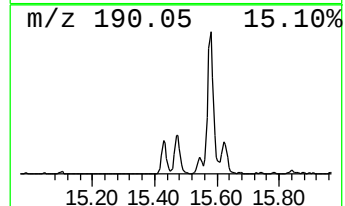
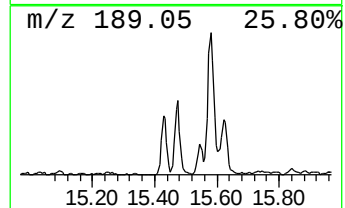
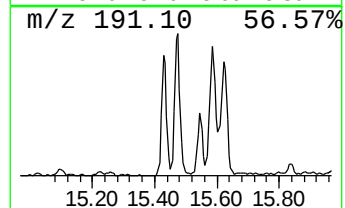
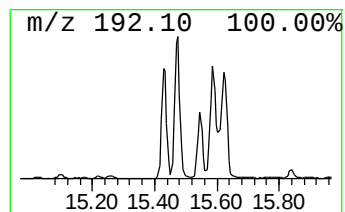
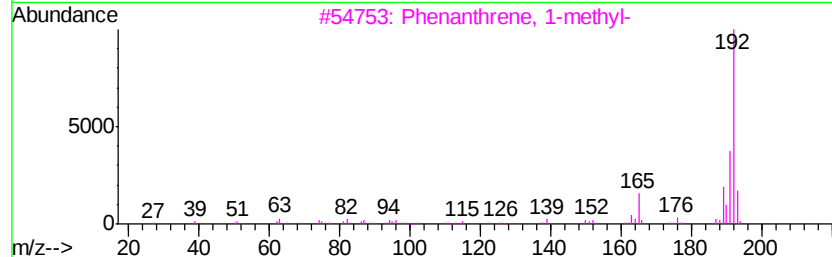
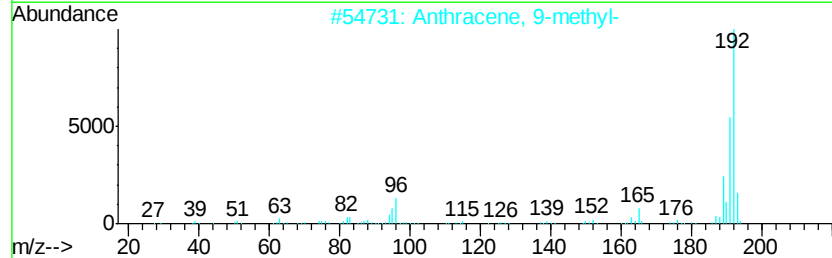
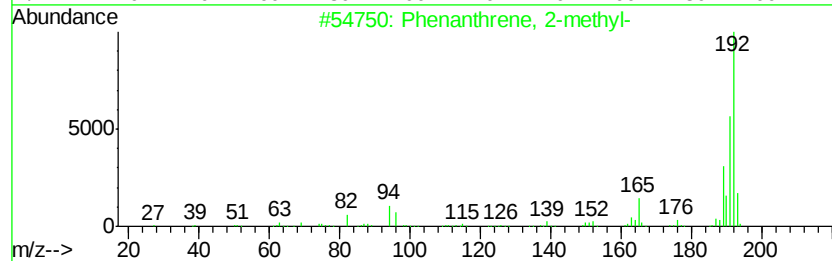
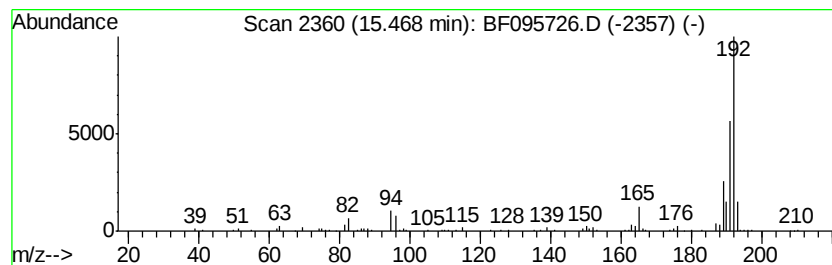
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.47 | 9.63 ng | 222308 | Phenanthrene-d10 | 14.63 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 98   |
| 2    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 95   |
| 3    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 95   |
| 4    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 95   |
| 5    |    |   | Phenanthrene, 3-methyl- | 192 | C15H12  | 000832-71-3 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

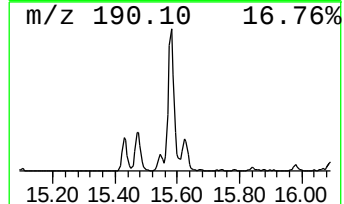
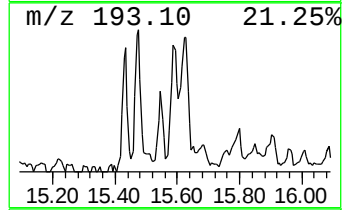
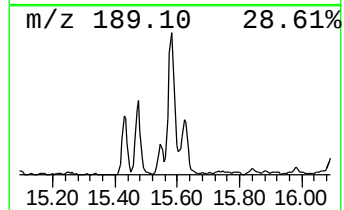
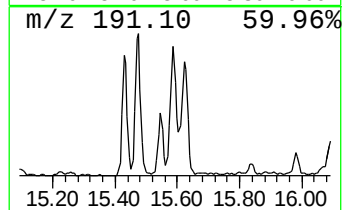
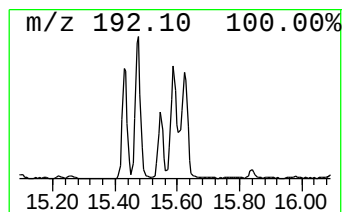
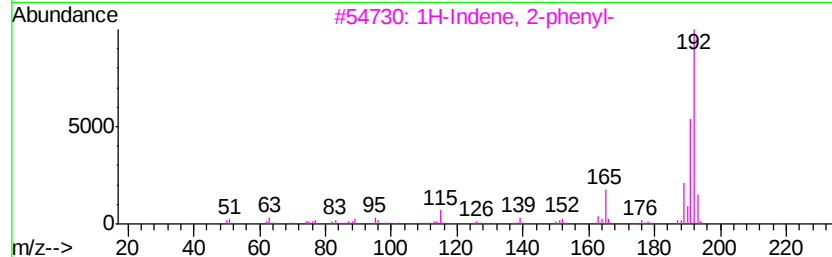
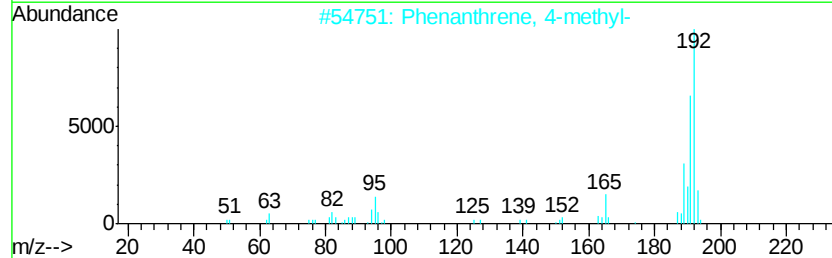
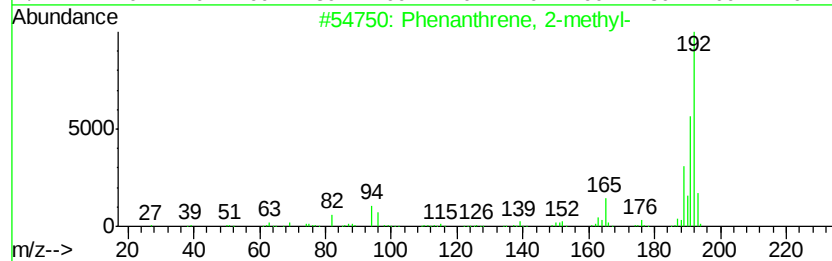
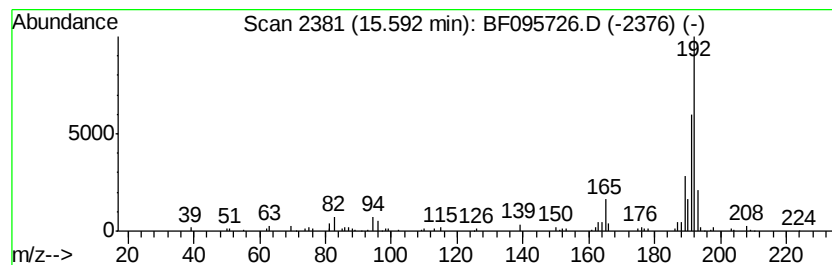
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Naphtho[2,3-b]norbornadiene Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.59 | 12.62 ng | 291281 | Phenanthrene-d10 | 14.63 |

| Hit# of | 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|---------|---|---------------------------------------|-----|---------|-------------|------|
| 1       |   | Phenanthrene, 2-methyl-               | 192 | C15H12  | 002531-84-2 | 70   |
| 2       |   | Phenanthrene, 4-methyl-               | 192 | C15H12  | 000832-64-4 | 64   |
| 3       |   | 1H-Indene, 2-phenyl-                  | 192 | C15H12  | 004505-48-0 | 64   |
| 4       |   | 1H-Cyclopropa[1]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 64   |
| 5       |   | Naphtho[2,3-b]norbornadiene           | 192 | C15H12  | 107426-38-0 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

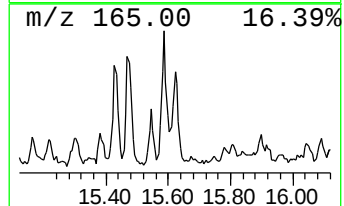
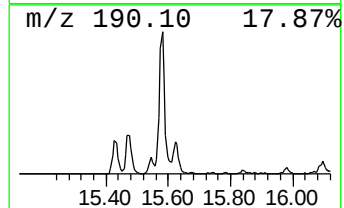
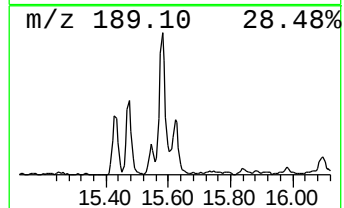
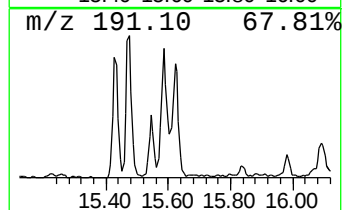
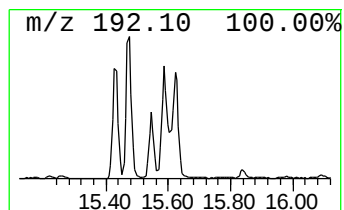
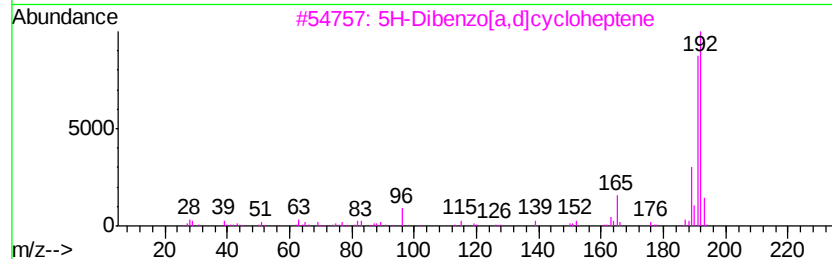
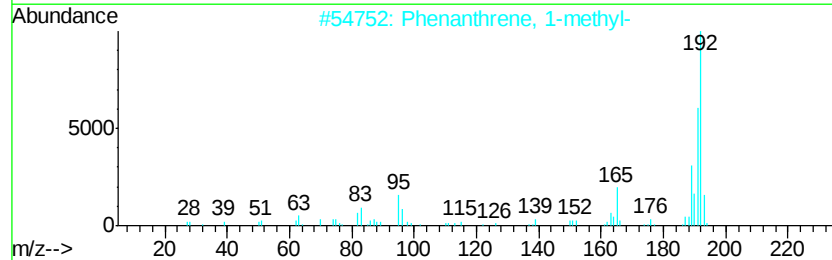
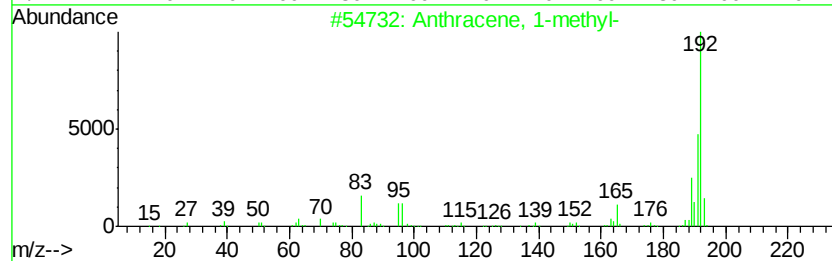
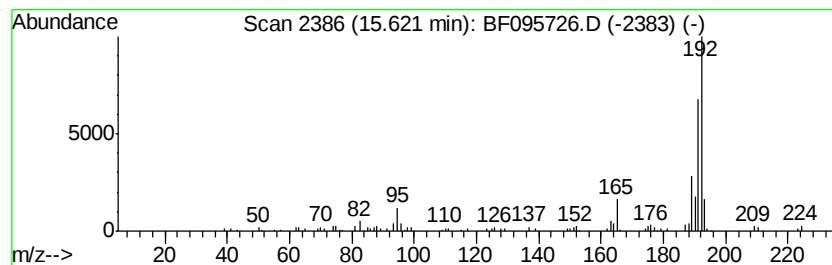
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Anthracene, 1-methyl- Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.62 | 12.32 ng | 284304 | Phenanthrene-d10 | 14.63 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|---------|---|-----------------------------|-----|---------|-------------|------|
| 1       |   | Anthracene, 1-methyl-       | 192 | C15H12  | 000610-48-0 | 95   |
| 2       |   | Phenanthrene, 1-methyl-     | 192 | C15H12  | 000832-69-9 | 90   |
| 3       |   | 5H-Dibenzo[a,d]cycloheptene | 192 | C15H12  | 000256-81-5 | 90   |
| 4       |   | Phenanthrene, 4-methyl-     | 192 | C15H12  | 000832-64-4 | 90   |
| 5       |   | Anthracene, 2-methyl-       | 192 | C15H12  | 000613-12-7 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

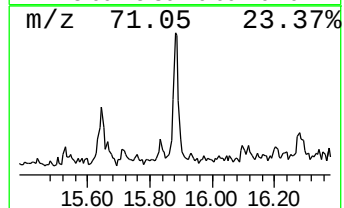
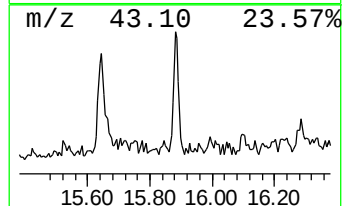
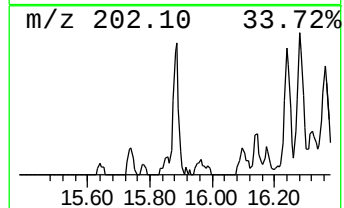
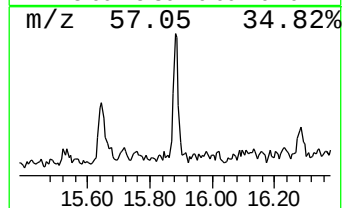
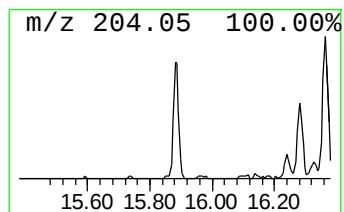
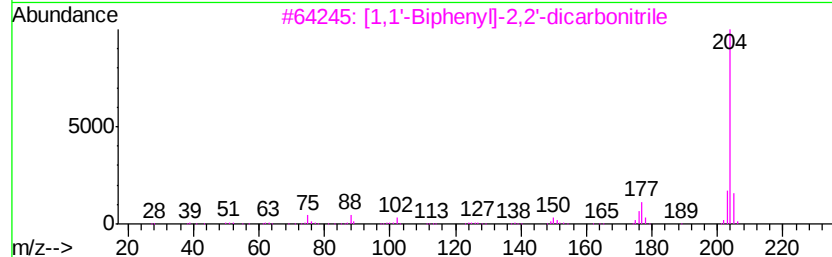
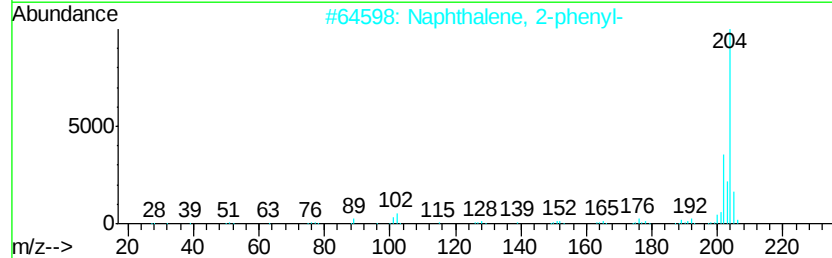
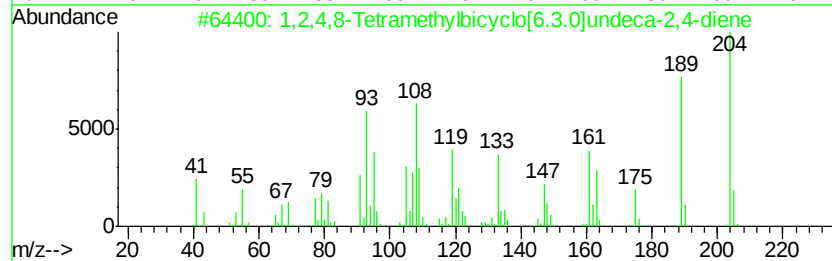
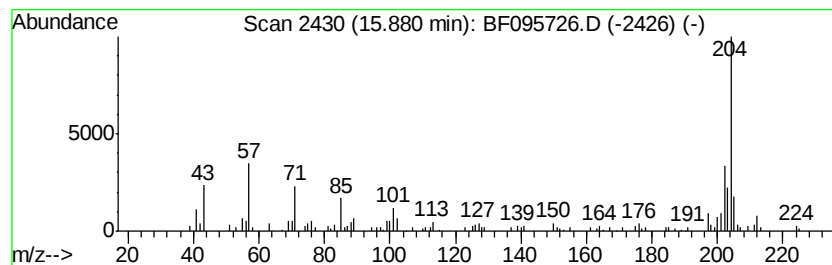
Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 1,2,4,8-Tetramethylbicyclo[...] Concentration Rank 11

| R.T.      | EstConc                              | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------------|--------|------------------|-------------|------|
| 15.88     | 7.09 ng                              | 163593 | Phenanthrene-d10 | 14.63       |      |
| Hit# of 5 | Tentative ID                         | MW     | MolForm          | CAS#        | Qual |
| 1         | 1,2,4,8-Tetramethylbicyclo[6.3.0...] | 204    | C15H24           | 137235-51-9 | 86   |
| 2         | Naphthalene, 2-phenyl-               | 204    | C16H12           | 000612-94-2 | 60   |
| 3         | [1,1'-Biphenyl]-2,2'-dicarbonitrile  | 204    | C14H8N2          | 004341-02-0 | 49   |
| 4         | [1,1'-Biphenyl]-4,4'-dicarbonitrile  | 204    | C14H8N2          | 001591-30-6 | 49   |
| 5         | Acephenanthrylene, 4,5-dihydro-      | 204    | C16H12           | 006232-48-0 | 45   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

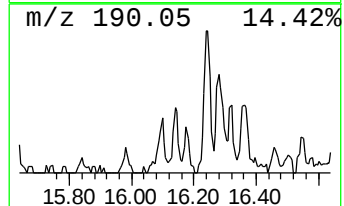
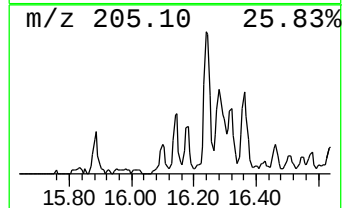
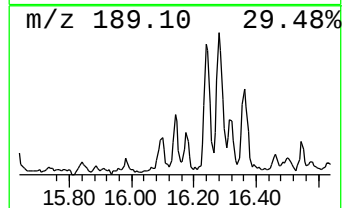
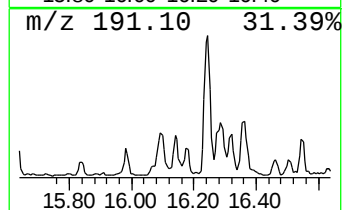
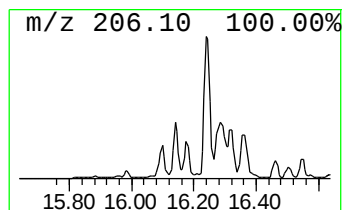
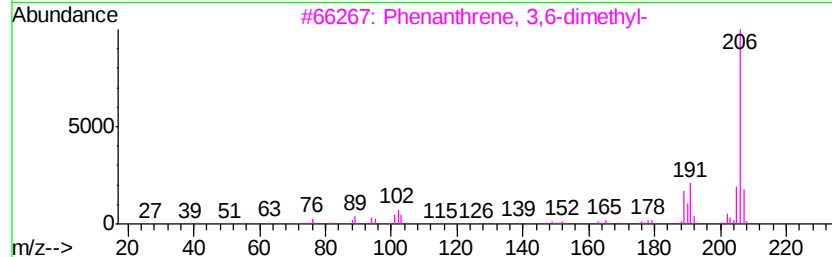
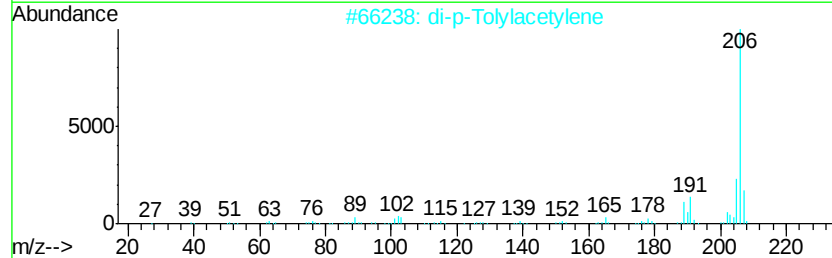
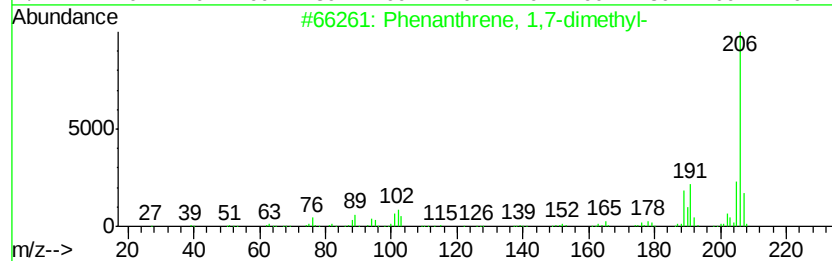
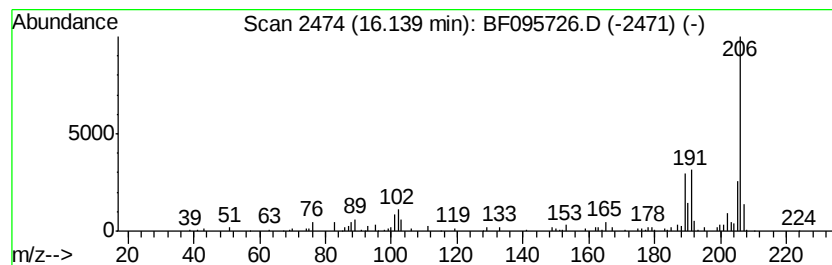
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Phenanthrene, 1,7-dimethyl- Concentration Rank 19

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.14 | 4.00 ng | 92247 | Phenanthrene-d10 | 14.63 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 94   |
| 2    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 91   |
| 3    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 91   |
| 4    |    |   | Phenanthrene, 2,7-dimethyl- | 206 | C16H14  | 001576-69-8 | 87   |
| 5    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 76   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

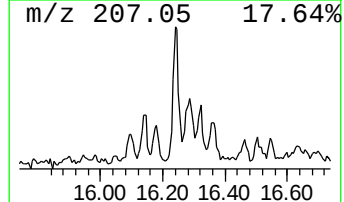
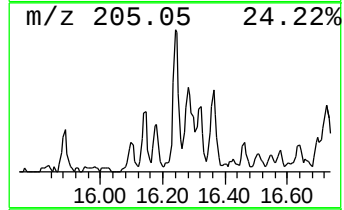
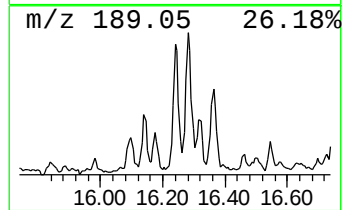
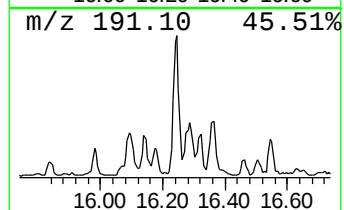
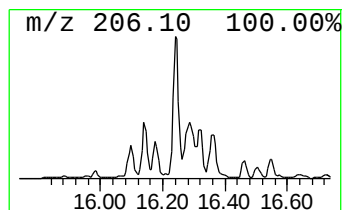
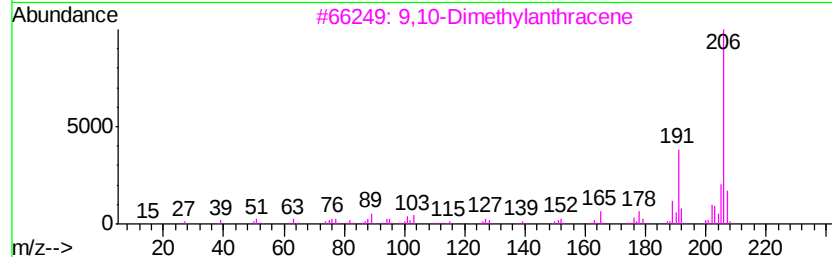
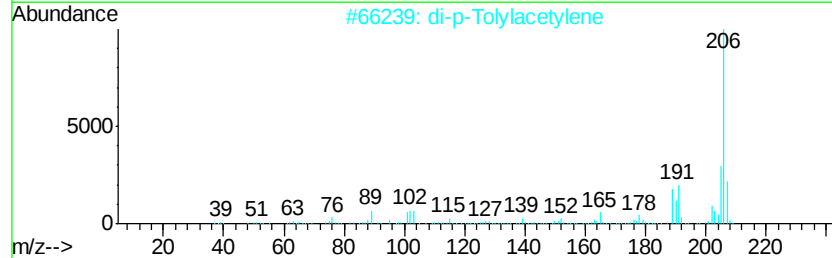
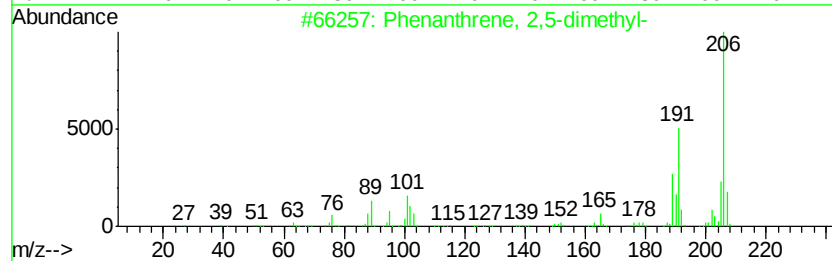
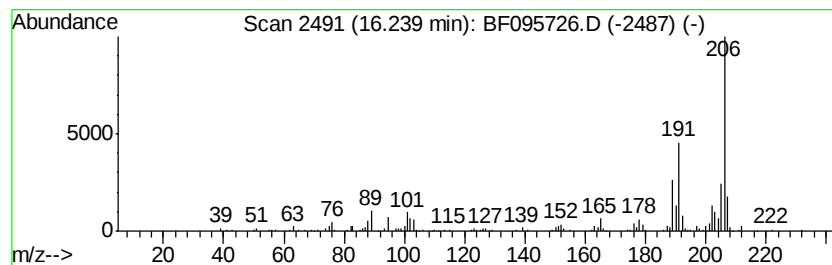
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Phenanthrene, 2,5-dimethyl- Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 16.24 | 12.87 ng | 297079 | Phenanthrene-d10 | 14.63 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 97   |
| 2    |    | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 96   |
| 3    |    | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 95   |
| 4    |    | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 94   |
| 5    |    | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

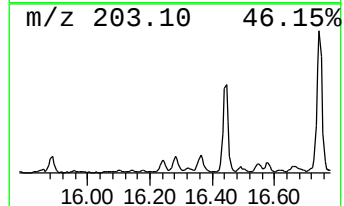
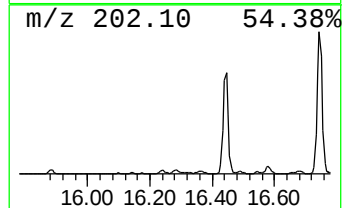
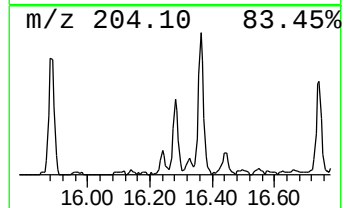
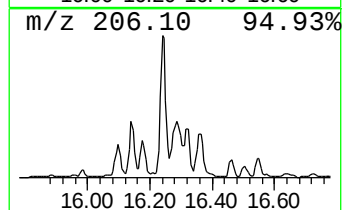
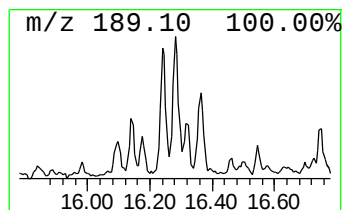
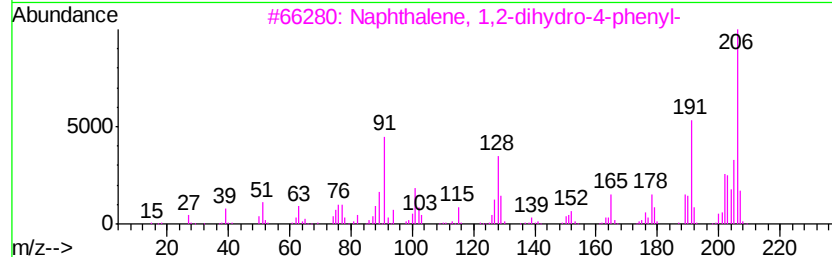
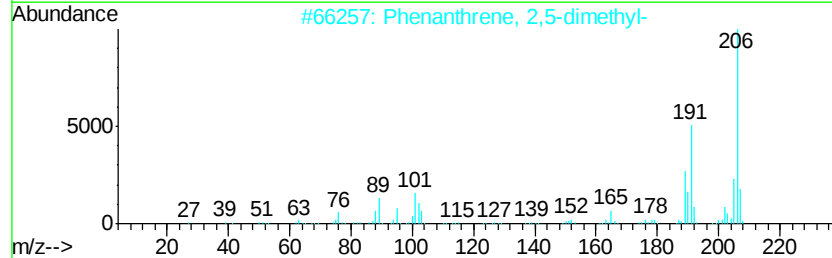
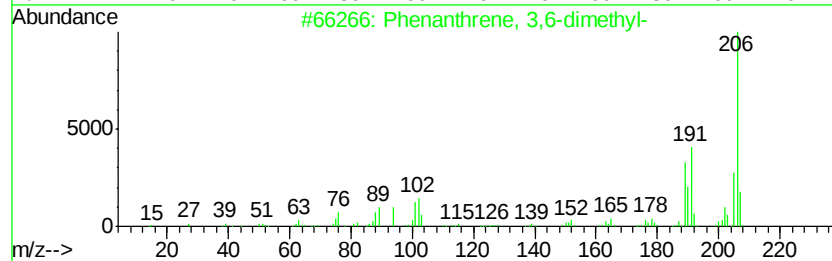
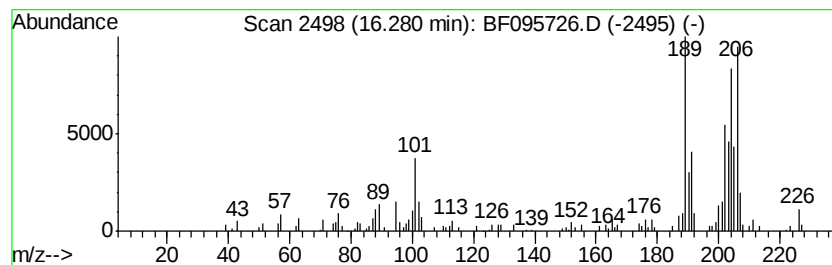
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Phenanthrene, 3,6-dimethyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.28 | 7.94 ng | 183391 | Phenanthrene-d10 | 14.63 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 3,6-dimethyl-         | 206 | C16H14  | 001576-67-6 | 83   |
| 2    |    | Phenanthrene, 2,5-dimethyl-         | 206 | C16H14  | 003674-66-6 | 42   |
| 3    |    | Naphthalene, 1,2-dihydro-4-phenyl-  | 206 | C16H14  | 007469-40-1 | 41   |
| 4    |    | Indeno[2,1-a]indene, 5,10-dihydro-  | 204 | C16H12  | 006543-29-9 | 41   |
| 5    |    | [14]Annulene, 1,6:8,13-bis(metha... | 206 | C16H14  | 085385-68-8 | 38   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

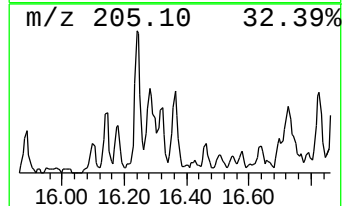
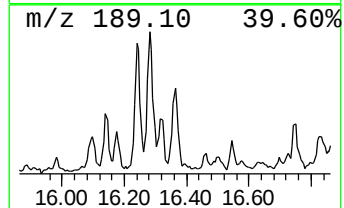
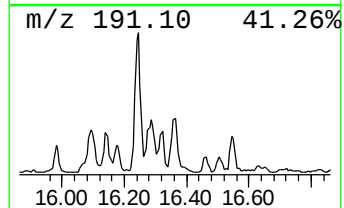
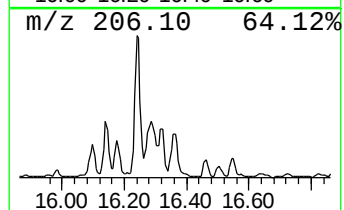
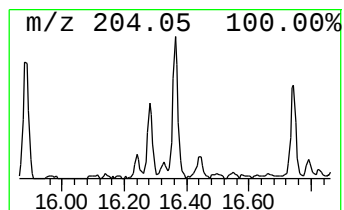
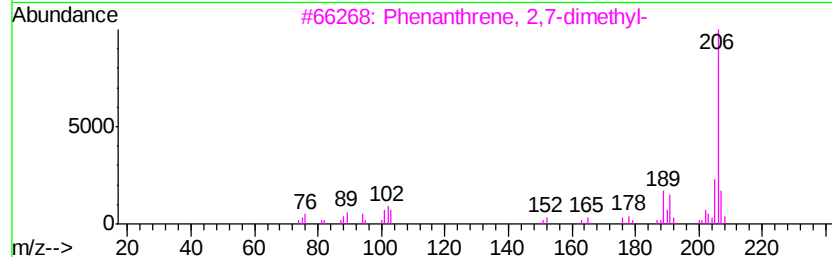
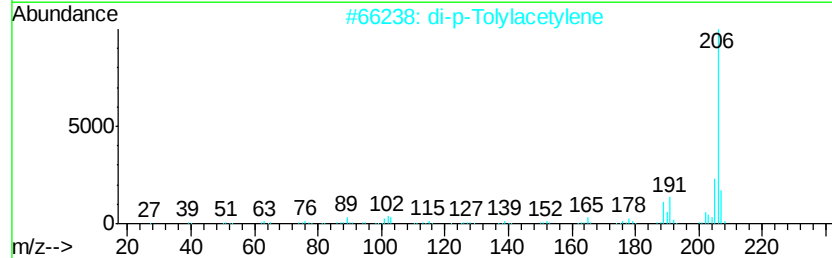
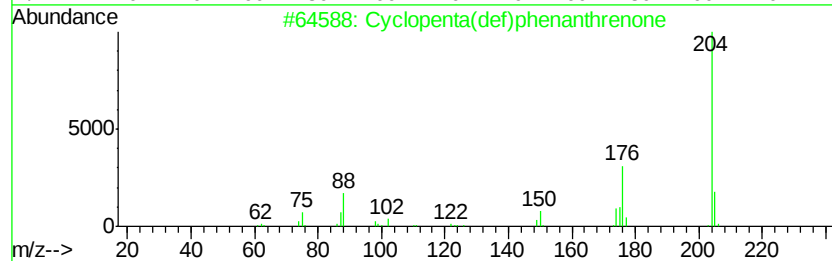
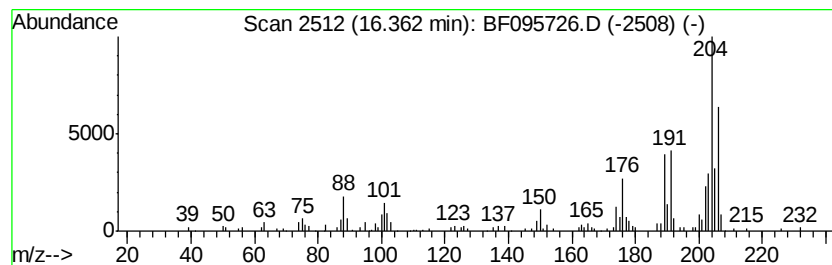
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Cyclopenta(def)phenanthrenone Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.36 | 7.57 ng | 174848 | Phenanthrene-d10 | 14.63 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrenone | 204 | C15H8O  | 005737-13-3 | 51   |
| 2    |    | di-p-Tolylacetylene           | 206 | C16H14  | 002789-88-0 | 49   |
| 3    |    | Phenanthrene, 2,7-dimethyl-   | 206 | C16H14  | 001576-69-8 | 45   |
| 4    |    | Phenanthrene, 3,6-dimethyl-   | 206 | C16H14  | 001576-67-6 | 45   |
| 5    |    | Phenanthrene, 1,7-dimethyl-   | 206 | C16H14  | 000483-87-4 | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

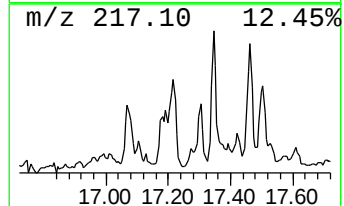
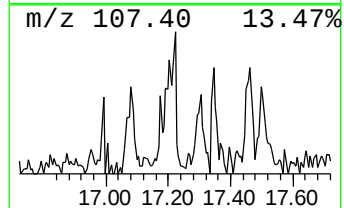
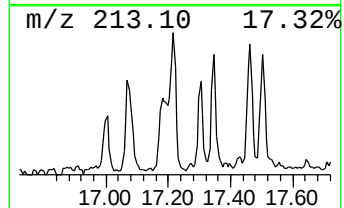
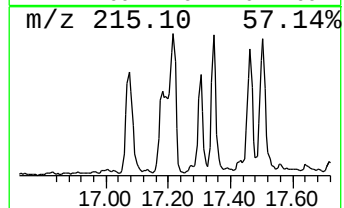
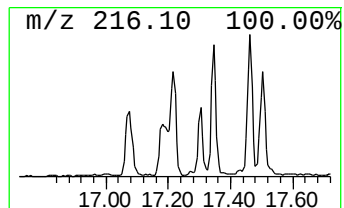
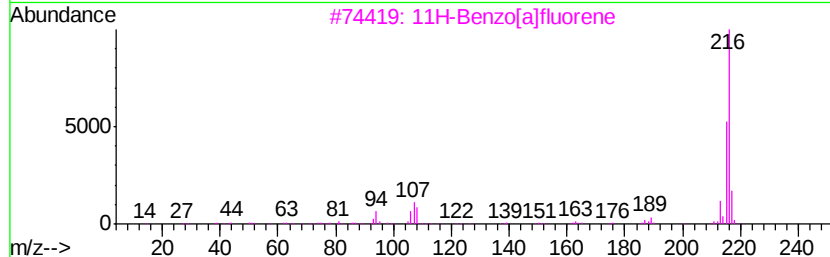
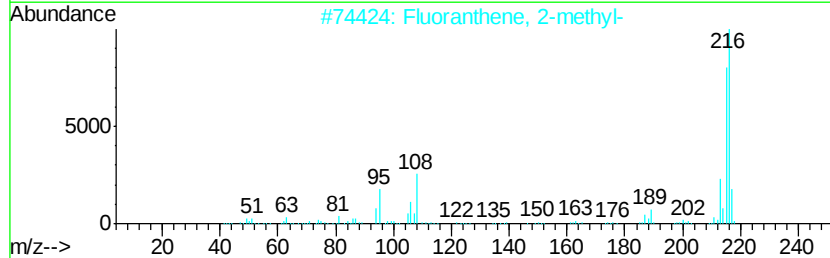
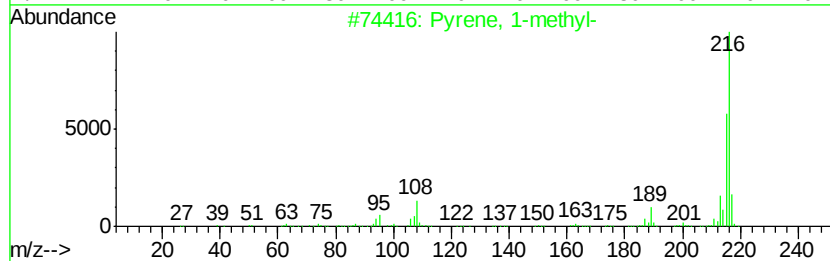
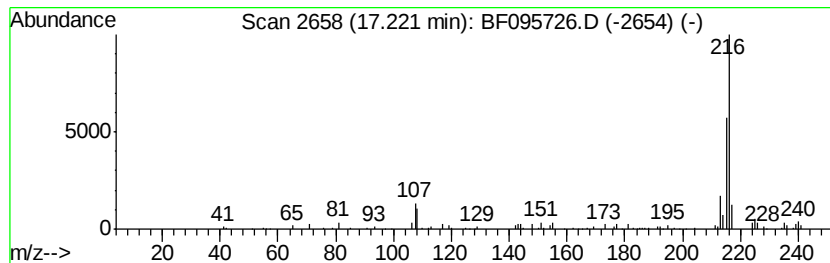
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Pyrene, 1-methyl- Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.22 | 5.09 ng | 229215 | Chrysene-d12     | 18.34 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 78   |
| 2    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 72   |
| 3    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 72   |
| 4    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 64   |
| 5    |    |   | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

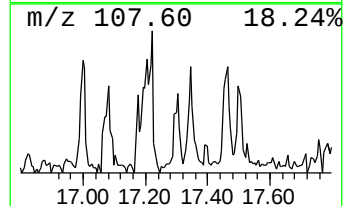
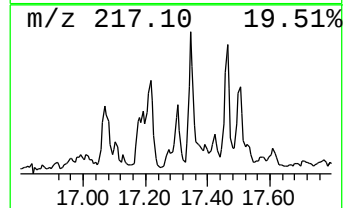
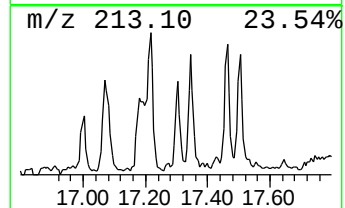
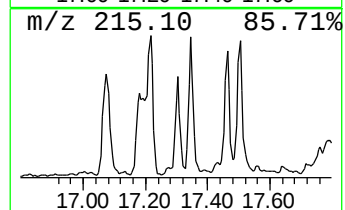
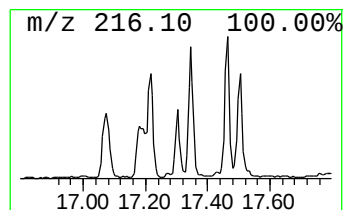
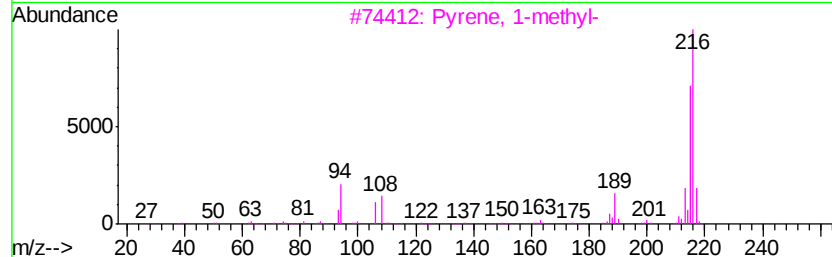
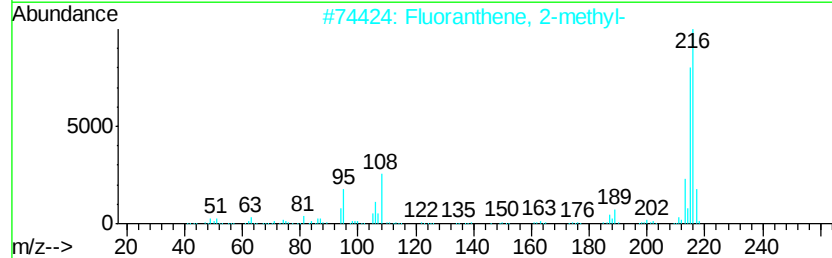
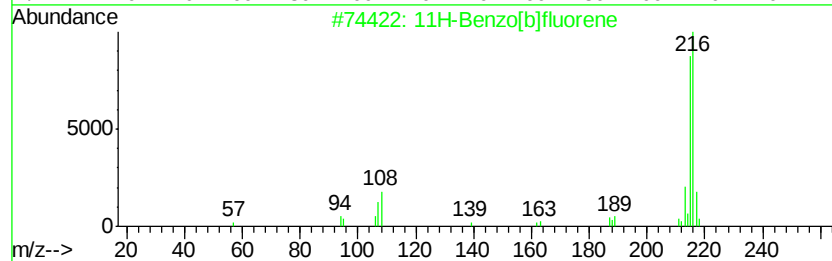
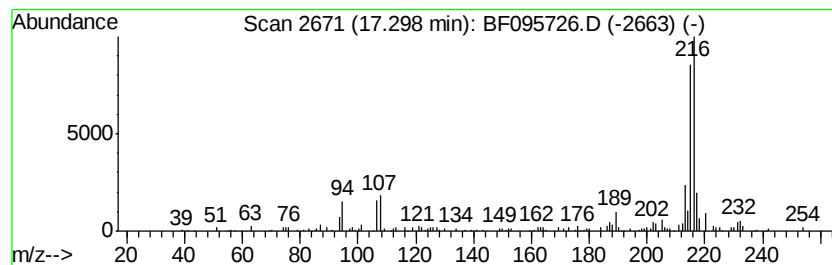
Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 18

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.    |             |      |
|---------|---------|-------------------------|------------------|---------|-------------|------|
| 17.30   | 4.67 ng | 210093                  | Chrysene-d12     | 18.34   |             |      |
| Hit# of | 5       | Tentative ID            | MW               | MolForm | CAS#        | Qual |
| 1       |         | 11H-Benzo[b]fluorene    | 216              | C17H12  | 000243-17-4 | 94   |
| 2       |         | Fluoranthene, 2-methyl- | 216              | C17H12  | 033543-31-6 | 93   |
| 3       |         | Pyrene, 1-methyl-       | 216              | C17H12  | 002381-21-7 | 87   |
| 4       |         | 11H-Benzo[a]fluorene    | 216              | C17H12  | 000238-84-6 | 81   |
| 5       |         | 7H-Benzo[c]fluorene     | 216              | C17H12  | 000205-12-9 | 81   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

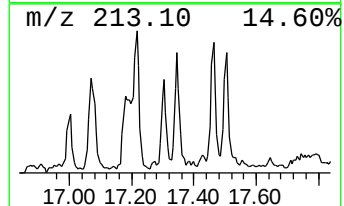
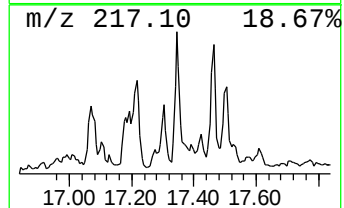
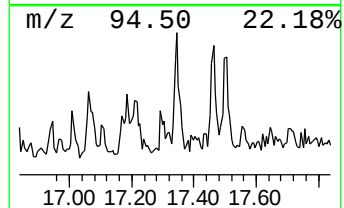
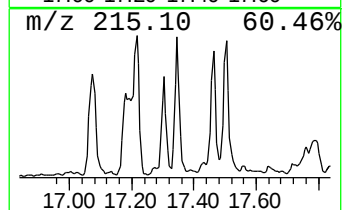
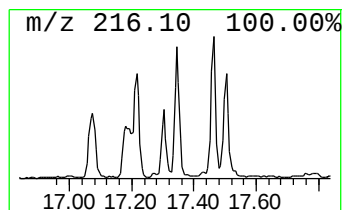
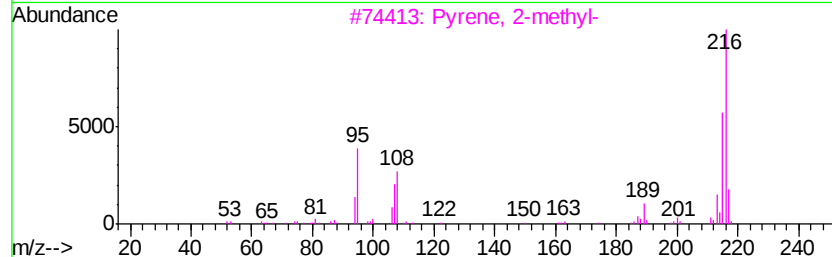
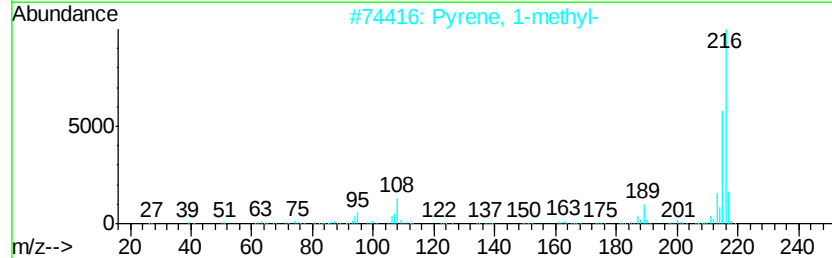
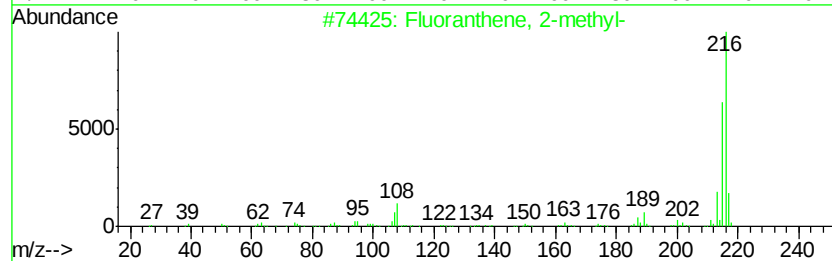
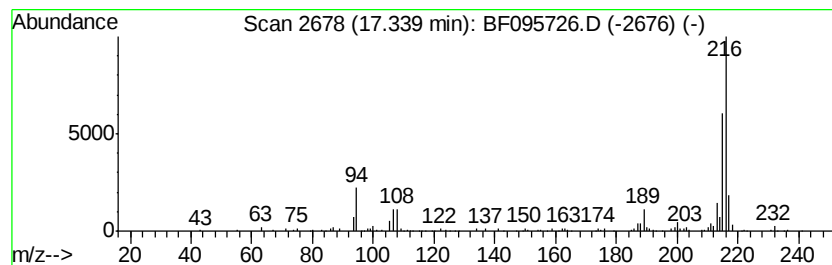
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Fluoranthene, 2-methyl- Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.34 | 5.85 ng | 263349 | Chrysene-d12     | 18.34 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 95   |
| 2    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 94   |
| 3    |    |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 93   |
| 4    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 90   |
| 5    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

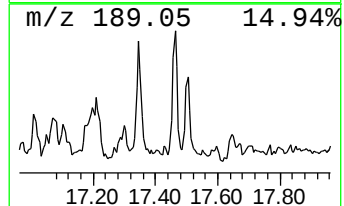
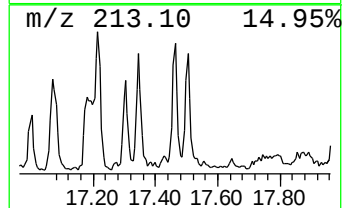
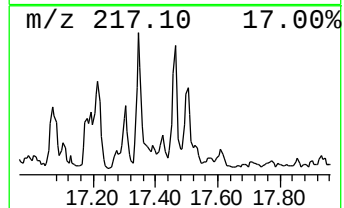
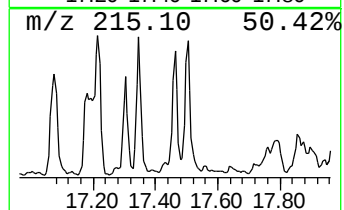
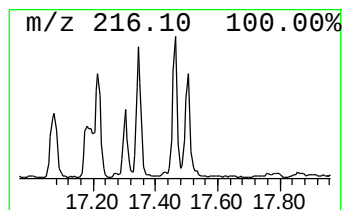
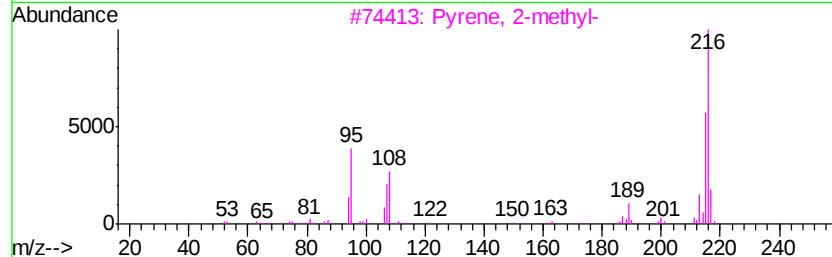
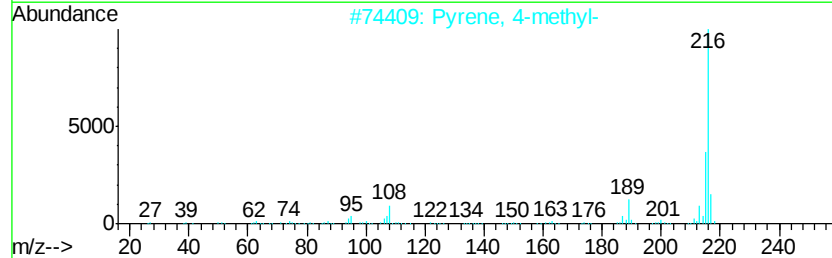
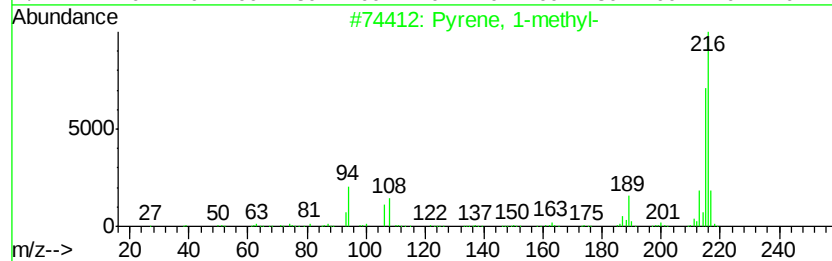
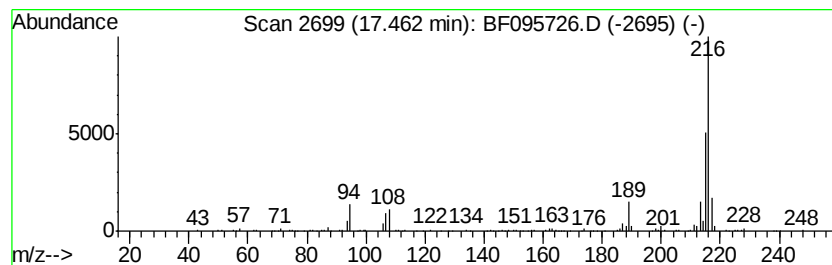
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Pyrene, 4-methyl- Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.46 | 3.98 ng | 179176 | Chrysene-d12     | 18.34 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 97   |
| 2    |    |   | Pyrene, 4-methyl-       | 216 | C17H12  | 003353-12-6 | 95   |
| 3    |    |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 95   |
| 4    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 91   |
| 5    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

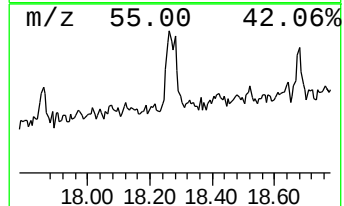
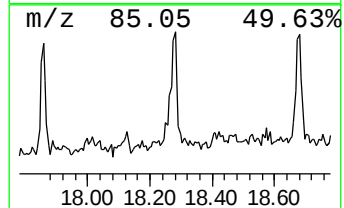
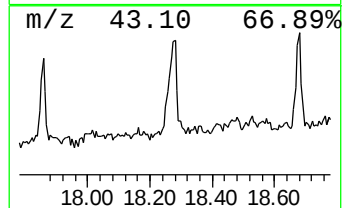
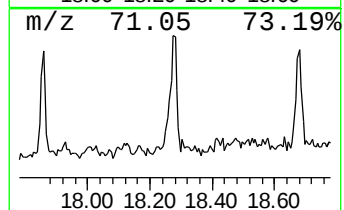
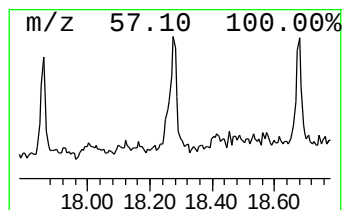
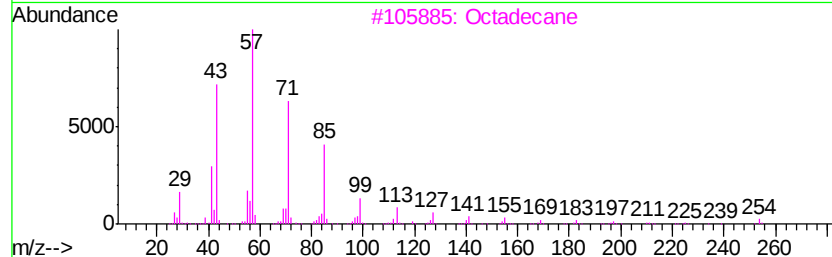
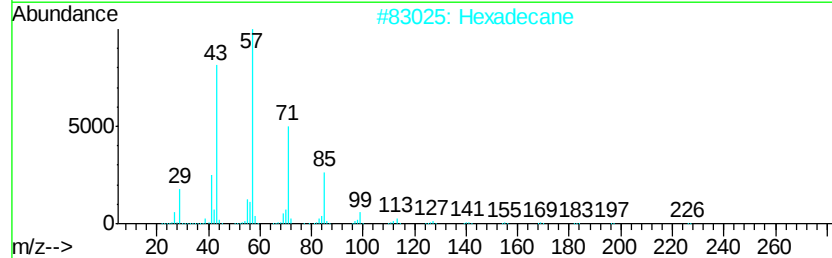
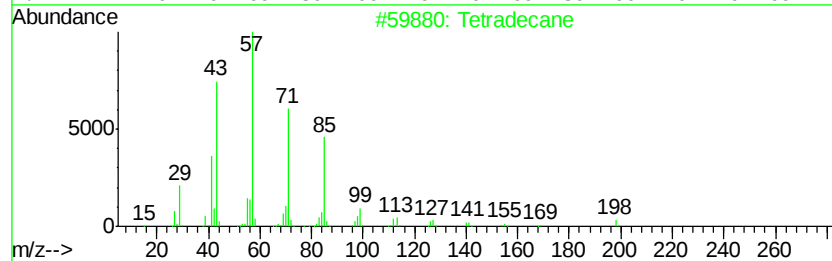
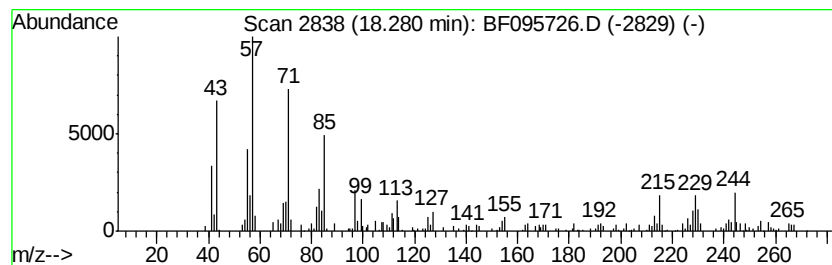
Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 Tetradecane Concentration Rank 17

| R.T.    | EstConc | Area           | Relative to ISTD | R.T.           |
|---------|---------|----------------|------------------|----------------|
| 18.28   | 4.79 ng | 215746         | Chrysene-d12     | 18.34          |
| Hit# of | 5       | Tentative ID   | MW MolForm       | CAS# Qual      |
| 1       |         | Tetradecane    | 198 C14H30       | 000629-59-4 90 |
| 2       |         | Hexadecane     | 226 C16H34       | 000544-76-3 83 |
| 3       |         | Octadecane     | 254 C18H38       | 000593-45-3 83 |
| 4       |         | Hentriacontane | 437 C31H64       | 000630-04-6 64 |
| 5       |         | Heptadecane    | 240 C17H36       | 000629-78-7 64 |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
Data File : BF095726.D  
Acq On : 7 Jun 2017 00:16  
Operator : SJ/MA  
Sample : I3482-07  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

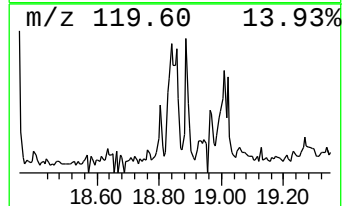
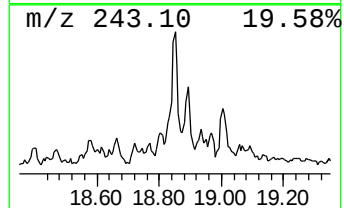
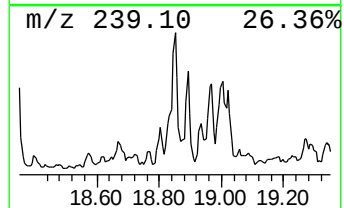
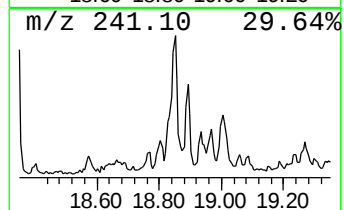
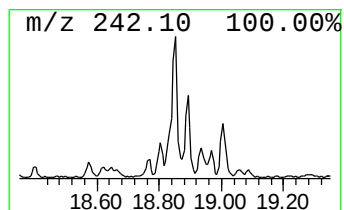
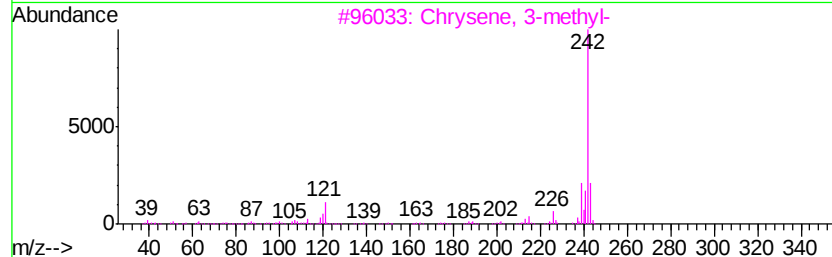
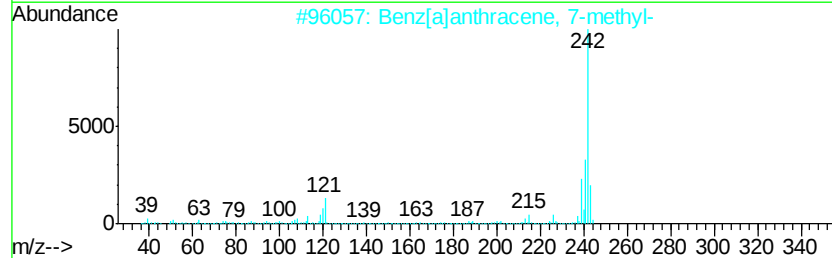
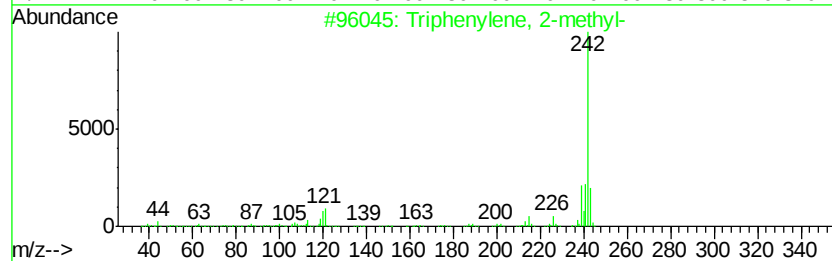
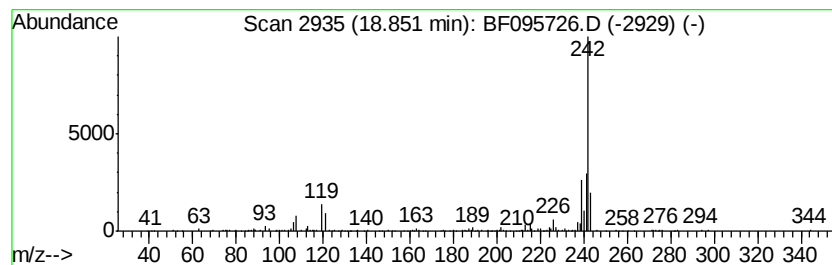
Instrument :  
BNA\_F  
ClientSampleId :  
CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 19 Triphenylene, 2-methyl- Concentration Rank 14

| R.T.    | EstConc | Area                         | Relative to ISTD |         | R.T.        |      |
|---------|---------|------------------------------|------------------|---------|-------------|------|
| 18.85   | 5.34 ng | 240310                       | Chrysene-d12     |         | 18.34       |      |
| Hit# of | 5       | Tentative ID                 | MW               | MolForm | CAS#        | Qual |
| 1       |         | Triphenylene, 2-methyl-      | 242              | C19H14  | 001705-84-6 | 95   |
| 2       |         | Benz[a]anthracene, 7-methyl- | 242              | C19H14  | 002541-69-7 | 94   |
| 3       |         | Chrysene, 3-methyl-          | 242              | C19H14  | 003351-31-3 | 94   |
| 4       |         | Benz[a]anthracene, 8-methyl- | 242              | C19H14  | 002381-31-9 | 92   |
| 5       |         | Chrysene, 5-methyl-          | 242              | C19H14  | 003697-24-3 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060617\  
 Data File : BF095726.D  
 Acq On : 7 Jun 2017 00:16  
 Operator : SJ/MA  
 Sample : I3482-07  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CL-01-060517-C

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060517.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| 2-Pentanone, 4-hy... | 4.59  | 5.9 ng  |       | 126800   | 1                     | 7.38  | 429972 | 20.0 |
| unknown7.03          | 7.03  | 76.0 ng |       | 1633470  | 1                     | 7.38  | 429972 | 20.0 |
| Dodecane             | 13.87 | 5.3 ng  |       | 122367   | 4                     | 14.63 | 461653 | 20.0 |
| Phenanthrene, 1-m... | 15.43 | 8.2 ng  |       | 189918   | 4                     | 14.63 | 461653 | 20.0 |
| Phenanthrene, 2-m... | 15.47 | 9.6 ng  |       | 222308   | 4                     | 14.63 | 461653 | 20.0 |
| Naphtho[2,3-b]nor... | 15.59 | 12.6 ng |       | 291281   | 4                     | 14.63 | 461653 | 20.0 |
| Anthracene, 1-met... | 15.62 | 12.3 ng |       | 284304   | 4                     | 14.63 | 461653 | 20.0 |
| 1,2,4,8-Tetrameth... | 15.88 | 7.1 ng  |       | 163593   | 4                     | 14.63 | 461653 | 20.0 |
| Phenanthrene, 1,7... | 16.14 | 4.0 ng  |       | 92247    | 4                     | 14.63 | 461653 | 20.0 |
| Phenanthrene, 2,5... | 16.24 | 12.9 ng |       | 297079   | 4                     | 14.63 | 461653 | 20.0 |
| Phenanthrene, 3,6... | 16.28 | 7.9 ng  |       | 183391   | 4                     | 14.63 | 461653 | 20.0 |
| Cyclopenta(def)ph... | 16.36 | 7.6 ng  |       | 174848   | 4                     | 14.63 | 461653 | 20.0 |
| Pyrene, 1-methyl-    | 17.22 | 5.1 ng  |       | 229215   | 5                     | 18.34 | 900642 | 20.0 |
| 11H-Benzo[b]fluorene | 17.30 | 4.7 ng  |       | 210093   | 5                     | 18.34 | 900642 | 20.0 |
| Fluoranthene, 2-m... | 17.34 | 5.8 ng  |       | 263349   | 5                     | 18.34 | 900642 | 20.0 |
| Pyrene, 4-methyl-    | 17.46 | 4.0 ng  |       | 179176   | 5                     | 18.34 | 900642 | 20.0 |
| Tetradecane          | 18.28 | 4.8 ng  |       | 215746   | 5                     | 18.34 | 900642 | 20.0 |
| Triphenylene, 2-m... | 18.85 | 5.3 ng  |       | 240310   | 5                     | 18.34 | 900642 | 20.0 |