

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
 Data File : BF085443.D  
 Acq On : 14 Mar 2016 21:53  
 Operator : UM/SJ  
 Sample : H1722-10  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 W170TH-SB-22(0.5-2.0)

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-BF030316.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         | 5.133       | 247           | 249         | 257          | rVB      | 211867         | 306428        | 6.73%           | 0.460%        |
| 2         | 5.533       | 281           | 284         | 287          | rBV      | 3514903        | 3754074       | 82.43%          | 5.640%        |
| 3         | 6.562       | 371           | 374         | 376          | rBV      | 3450664        | 4554250       | 100.00%         | 6.842%        |
| 4         | 6.687       | 383           | 385         | 388          | rBV      | 3048074        | 4390354       | 96.40%          | 6.596%        |
| 5         | 6.927       | 403           | 406         | 408          | rBV      | 798990         | 1060751       | 23.29%          | 1.594%        |
| 6         | 7.076       | 417           | 419         | 422          | rVB      | 2831013        | 3274765       | 71.91%          | 4.920%        |
| 7         | 7.487       | 452           | 455         | 457          | rBV      | 2640181        | 2713646       | 59.58%          | 4.077%        |
| 8         | 7.579       | 460           | 463         | 465          | rVB      | 87933          | 108322        | 2.38%           | 0.163%        |
| 9         | 8.207       | 515           | 518         | 520          | rBV      | 1776203        | 1530915       | 33.62%          | 2.300%        |
| 10        | 9.282       | 609           | 612         | 615          | rBV      | 3448714        | 4377084       | 96.11%          | 6.576%        |
| 11        | 9.625       | 639           | 642         | 645          | rBV      | 93409          | 161011        | 3.54%           | 0.242%        |
| 12        | 9.682       | 645           | 647         | 648          | rVV      | 747433         | 835216        | 18.34%          | 1.255%        |
| 13        | 9.739       | 650           | 652         | 657          | rVB      | 59361          | 72091         | 1.58%           | 0.108%        |
| 14        | 9.819       | 657           | 659         | 661          | rBV      | 96650          | 107826        | 2.37%           | 0.162%        |
| 15        | 9.888       | 661           | 665         | 668          | rBV      | 112567         | 163740        | 3.60%           | 0.246%        |
| 16        | 9.968       | 669           | 672         | 673          | rVV      | 1570992        | 1562158       | 34.30%          | 2.347%        |
| 17        | 9.990       | 673           | 674         | 676          | rVB      | 108112         | 124557        | 2.73%           | 0.187%        |
| 18        | 10.116      | 683           | 685         | 687          | rBV3     | 39756          | 76971         | 1.69%           | 0.116%        |
| 19        | 10.162      | 687           | 689         | 692          | rVV3     | 125184         | 151732        | 3.33%           | 0.228%        |
| 20        | 10.208      | 692           | 693         | 696          | rVB2     | 77825          | 72098         | 1.58%           | 0.108%        |
| 21        | 10.276      | 698           | 699         | 701          | rBV      | 64961          | 67816         | 1.49%           | 0.102%        |
| 22        | 10.368      | 705           | 707         | 708          | rBV      | 85047          | 103507        | 2.27%           | 0.156%        |
| 23        | 10.390      | 708           | 709         | 711          | rVB      | 248484         | 193287        | 4.24%           | 0.290%        |
| 24        | 10.505      | 718           | 719         | 722          | rVB      | 195535         | 219292        | 4.82%           | 0.329%        |
| 25        | 10.573      | 722           | 725         | 726          | rBV2     | 55724          | 78272         | 1.72%           | 0.118%        |
| 26        | 10.619      | 726           | 729         | 732          | rVV3     | 68777          | 138354        | 3.04%           | 0.208%        |
| 27        | 10.665      | 732           | 733         | 735          | rVB      | 107520         | 83939         | 1.84%           | 0.126%        |
| 28        | 10.756      | 738           | 741         | 743          | rBV      | 2784473        | 2903731       | 63.76%          | 4.363%        |
| 29        | 10.871      | 749           | 751         | 755          | rVB      | 210425         | 329243        | 7.23%           | 0.495%        |
| 30        | 11.053      | 765           | 767         | 769          | rBV3     | 122004         | 188623        | 4.14%           | 0.283%        |
| 31        | 11.191      | 776           | 779         | 783          | rVV4     | 56607          | 165522        | 3.63%           | 0.249%        |
| 32        | 11.316      | 786           | 790         | 791          | rBV2     | 145016         | 213814        | 4.69%           | 0.321%        |
| 33        | 11.339      | 791           | 792         | 795          | rVB      | 117549         | 170195        | 3.74%           | 0.256%        |
| 34        | 11.453      | 799           | 802         | 803          | rBV      | 1349393        | 1891282       | 41.53%          | 2.842%        |

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 W170TH-SB-22(0.5-2.0)

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-BF030316.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 11.476 | 803  | 804  | 806  | rVV  | 1867766 | 1326981 | 29.14% | 1.994% |
| 36 | 11.522 | 806  | 808  | 810  | rVV  | 662382  | 553628  | 12.16% | 0.832% |
| 37 | 11.682 | 818  | 822  | 825  | rBV2 | 50698   | 144337  | 3.17%  | 0.217% |
| 38 | 11.739 | 825  | 827  | 829  | rVB  | 119195  | 105931  | 2.33%  | 0.159% |
| 39 | 11.773 | 829  | 830  | 833  | rVB2 | 71456   | 94795   | 2.08%  | 0.142% |
| 40 | 11.945 | 843  | 845  | 846  | rBV  | 259644  | 337038  | 7.40%  | 0.506% |
| 41 | 11.979 | 846  | 848  | 850  | rVB  | 1013209 | 908126  | 19.94% | 1.364% |
| 42 | 12.025 | 850  | 852  | 853  | rBV  | 209992  | 164544  | 3.61%  | 0.247% |
| 43 | 12.059 | 853  | 855  | 859  | rVB  | 663941  | 1009000 | 22.16% | 1.516% |
| 44 | 12.151 | 859  | 863  | 864  | rBV2 | 64517   | 96862   | 2.13%  | 0.146% |
| 45 | 12.242 | 869  | 871  | 872  | rBV  | 240258  | 202137  | 4.44%  | 0.304% |
| 46 | 12.391 | 881  | 884  | 886  | rBV2 | 75902   | 103949  | 2.28%  | 0.156% |
| 47 | 12.425 | 886  | 887  | 888  | rVB  | 112364  | 77082   | 1.69%  | 0.116% |
| 48 | 12.505 | 891  | 894  | 895  | rBV  | 200647  | 312829  | 6.87%  | 0.470% |
| 49 | 12.539 | 895  | 897  | 900  | rVV  | 164703  | 232191  | 5.10%  | 0.349% |
| 50 | 12.585 | 900  | 901  | 906  | rVB3 | 176145  | 201568  | 4.43%  | 0.303% |
| 51 | 12.665 | 906  | 908  | 910  | rBV  | 3003689 | 2714986 | 59.61% | 4.079% |
| 52 | 12.722 | 910  | 913  | 915  | rVV2 | 64559   | 129028  | 2.83%  | 0.194% |
| 53 | 12.756 | 915  | 916  | 920  | rVV3 | 259333  | 313213  | 6.88%  | 0.471% |
| 54 | 12.814 | 920  | 921  | 923  | rVV  | 113010  | 100604  | 2.21%  | 0.151% |
| 55 | 12.894 | 925  | 928  | 930  | rVV  | 2461993 | 2702171 | 59.33% | 4.060% |
| 56 | 12.939 | 930  | 932  | 934  | rVB2 | 120466  | 154969  | 3.40%  | 0.233% |
| 57 | 13.031 | 938  | 940  | 943  | rVB  | 2457549 | 3390242 | 74.44% | 5.094% |
| 58 | 13.111 | 943  | 947  | 951  | rBV3 | 189679  | 333027  | 7.31%  | 0.500% |
| 59 | 13.214 | 953  | 956  | 958  | rVB  | 352051  | 544595  | 11.96% | 0.818% |
| 60 | 13.282 | 960  | 962  | 964  | rVV  | 389304  | 351136  | 7.71%  | 0.528% |
| 61 | 13.316 | 964  | 965  | 969  | rVB2 | 217059  | 308376  | 6.77%  | 0.463% |
| 62 | 13.419 | 972  | 974  | 975  | rVB  | 101850  | 130517  | 2.87%  | 0.196% |
| 63 | 13.442 | 975  | 976  | 978  | rBV  | 158082  | 154887  | 3.40%  | 0.233% |
| 64 | 13.591 | 987  | 989  | 990  | rBV2 | 60377   | 80015   | 1.76%  | 0.120% |
| 65 | 13.728 | 997  | 1001 | 1003 | rBV2 | 90224   | 197681  | 4.34%  | 0.297% |
| 66 | 13.831 | 1008 | 1010 | 1012 | rBV2 | 145713  | 259452  | 5.70%  | 0.390% |
| 67 | 13.865 | 1012 | 1013 | 1014 | rVV  | 195722  | 157460  | 3.46%  | 0.237% |
| 68 | 13.888 | 1014 | 1015 | 1017 | rVV2 | 149151  | 212055  | 4.66%  | 0.319% |
| 69 | 13.922 | 1017 | 1018 | 1022 | rVV2 | 710402  | 838141  | 18.40% | 1.259% |
| 70 | 14.002 | 1023 | 1025 | 1028 | rVV  | 52211   | 81459   | 1.79%  | 0.122% |
| 71 | 14.082 | 1029 | 1032 | 1034 | rVV2 | 1473828 | 2408509 | 52.88% | 3.619% |

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 ClientSampleId :  
 W170TH-SB-22(0.5-2.0)

Integration Parameters: rteint.p  
 Integrator: RTE  
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 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-BF030316.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |        |         |        |        |
|-----|--------|------|------|------|------|--------|---------|--------|--------|
| 72  | 14.117 | 1034 | 1035 | 1037 | rVV  | 820286 | 646843  | 14.20% | 0.972% |
| 73  | 14.197 | 1038 | 1042 | 1043 | rVV2 | 137355 | 211377  | 4.64%  | 0.318% |
| 74  | 14.231 | 1043 | 1045 | 1048 | rVV2 | 74678  | 201336  | 4.42%  | 0.302% |
| 75  | 14.311 | 1050 | 1052 | 1054 | rVV2 | 84957  | 144040  | 3.16%  | 0.216% |
| 76  | 14.414 | 1058 | 1061 | 1062 | rBV2 | 46639  | 80360   | 1.76%  | 0.121% |
| 77  | 14.437 | 1062 | 1063 | 1064 | rVV  | 68684  | 69323   | 1.52%  | 0.104% |
| 78  | 14.471 | 1064 | 1066 | 1068 | rVV  | 231160 | 285902  | 6.28%  | 0.430% |
| 79  | 14.505 | 1068 | 1069 | 1071 | rVV  | 119459 | 112213  | 2.46%  | 0.169% |
| 80  | 14.562 | 1071 | 1074 | 1076 | rVV3 | 175073 | 294461  | 6.47%  | 0.442% |
| 81  | 14.619 | 1076 | 1079 | 1081 | rVV3 | 155702 | 279012  | 6.13%  | 0.419% |
| 82  | 14.665 | 1081 | 1083 | 1086 | rVB3 | 58151  | 98980   | 2.17%  | 0.149% |
| 83  | 14.779 | 1091 | 1093 | 1094 | rBV2 | 41583  | 70480   | 1.55%  | 0.106% |
| 84  | 14.837 | 1097 | 1098 | 1101 | rVB2 | 125777 | 155966  | 3.42%  | 0.234% |
| 85  | 15.122 | 1120 | 1123 | 1128 | rBV  | 720007 | 1464860 | 32.16% | 2.201% |
| 86  | 15.225 | 1130 | 1132 | 1134 | rVB  | 166803 | 177351  | 3.89%  | 0.266% |
| 87  | 15.328 | 1138 | 1141 | 1143 | rBV2 | 38234  | 107755  | 2.37%  | 0.162% |
| 88  | 15.420 | 1147 | 1149 | 1152 | rVB  | 426586 | 547408  | 12.02% | 0.822% |
| 89  | 15.477 | 1152 | 1154 | 1158 | rBV  | 551346 | 864799  | 18.99% | 1.299% |
| 90  | 15.545 | 1158 | 1160 | 1166 | rVB2 | 728755 | 1172013 | 25.73% | 1.761% |
| 91  | 16.060 | 1202 | 1205 | 1207 | rBV3 | 41390  | 70652   | 1.55%  | 0.106% |
| 92  | 16.242 | 1219 | 1221 | 1228 | rVB2 | 51238  | 129009  | 2.83%  | 0.194% |
| 93  | 16.985 | 1283 | 1286 | 1291 | rVB2 | 261949 | 550843  | 12.10% | 0.828% |
| 94  | 17.065 | 1291 | 1293 | 1299 | rVB2 | 33008  | 74498   | 1.64%  | 0.112% |
| 95  | 17.168 | 1299 | 1302 | 1304 | rBV  | 75647  | 156924  | 3.45%  | 0.236% |
| 96  | 17.225 | 1304 | 1307 | 1308 | rBV  | 38335  | 64464   | 1.42%  | 0.097% |
| 97  | 17.431 | 1321 | 1325 | 1333 | rVB  | 193248 | 479095  | 10.52% | 0.720% |
| 98  | 17.568 | 1333 | 1337 | 1342 | rVV5 | 50928  | 175146  | 3.85%  | 0.263% |
| 99  | 17.660 | 1342 | 1345 | 1352 | rVB2 | 70521  | 166401  | 3.65%  | 0.250% |
| 100 | 18.586 | 1424 | 1426 | 1434 | rVB9 | 38152  | 129199  | 2.84%  | 0.194% |

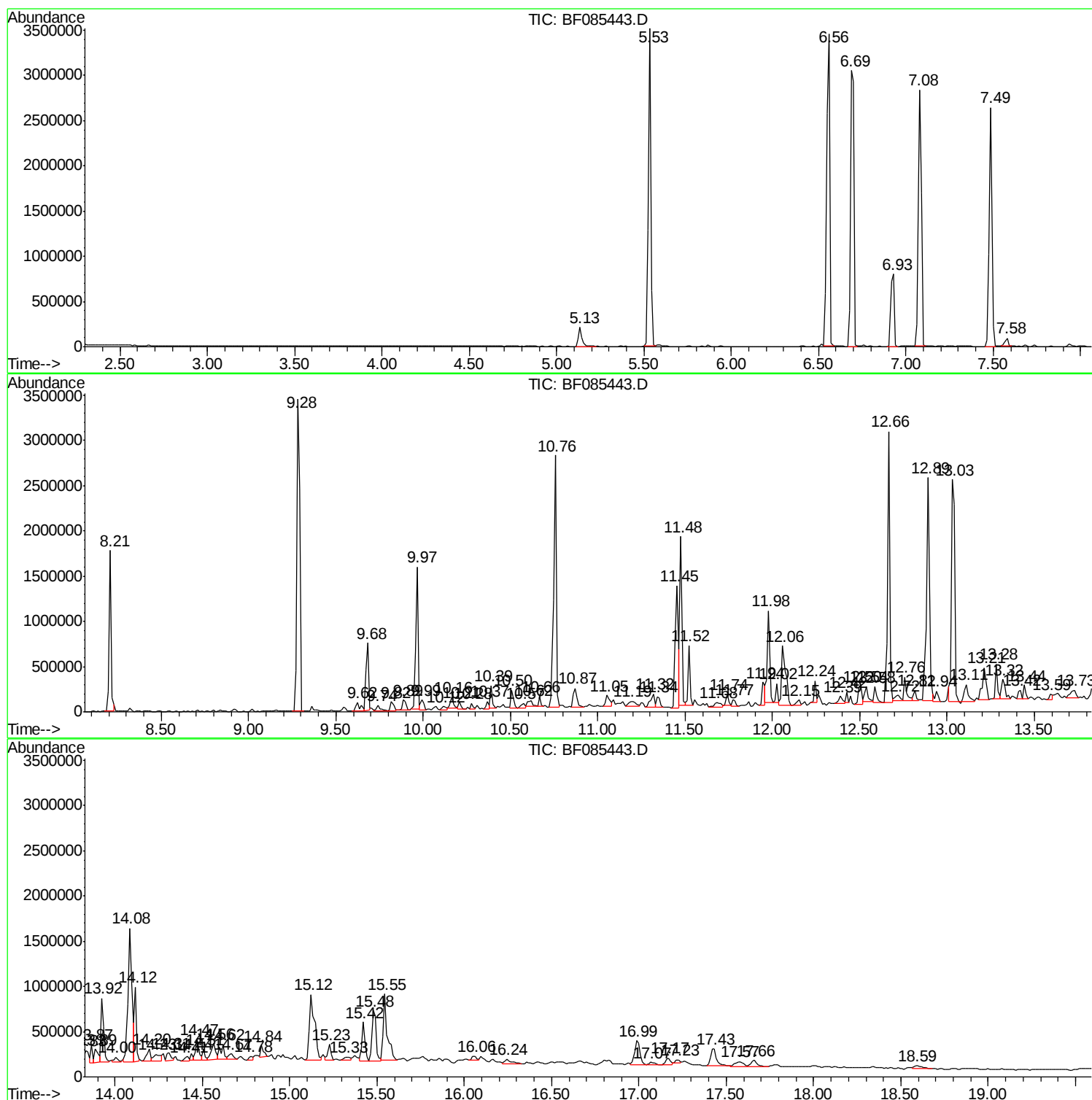
Sum of corrected areas: 66559097

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Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
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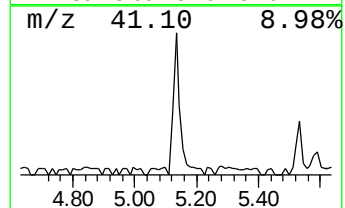
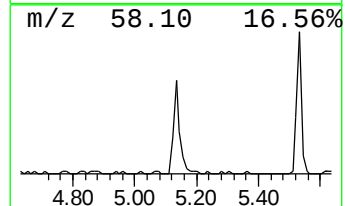
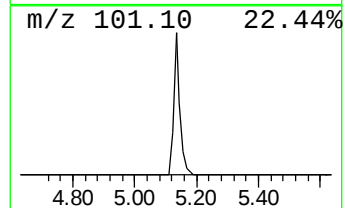
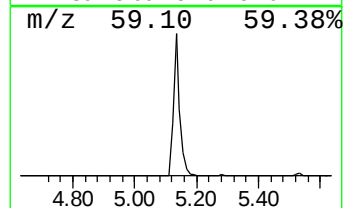
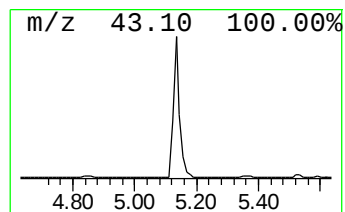
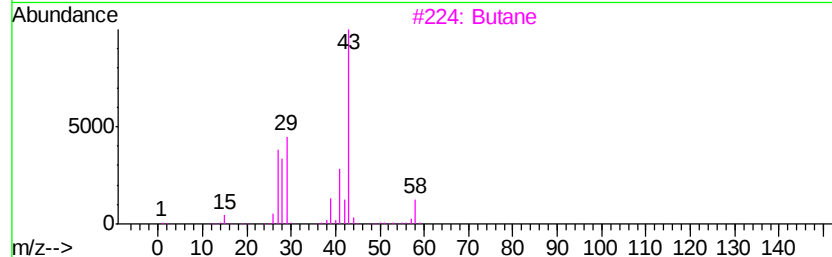
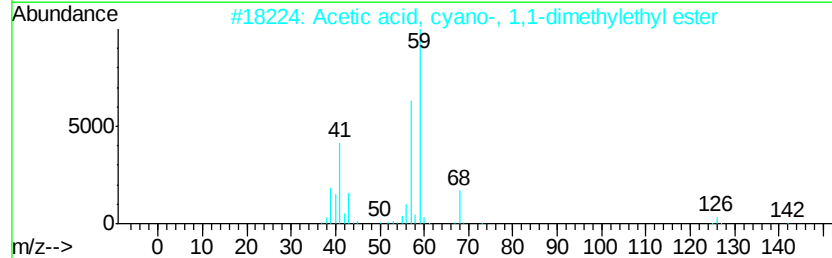
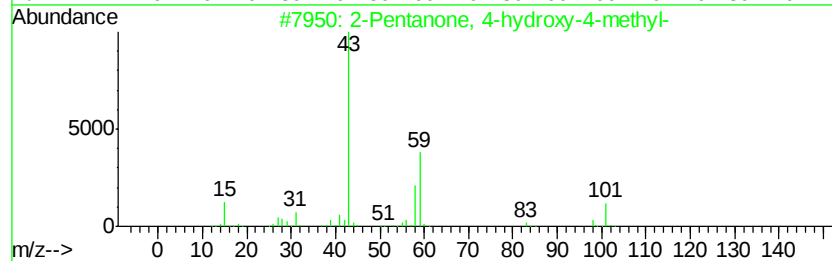
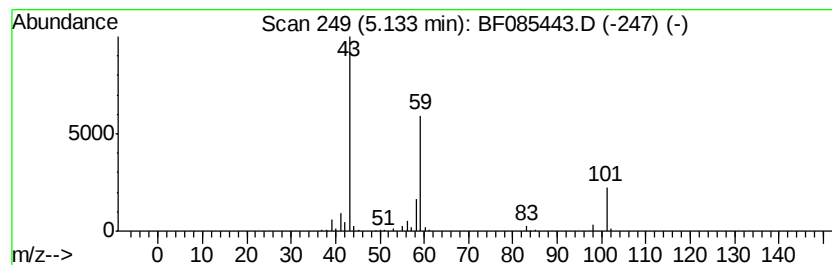
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.13 | 5.78 ng | 306428 | 1,4-Dichlorobenzene-d4 | 6.93 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2      | 000123-42-2 | 50      |      |      |
| 2    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2     | 001116-98-9 | 23      |      |      |
| 3    | Butane                              | 58  | C4H10        | 000106-97-8 | 9       |      |      |
| 4    | Morpholine, 4-methyl-               | 101 | C5H11NO      | 000109-02-4 | 9       |      |      |
| 5    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2      | 000057-71-6 | 9       |      |      |



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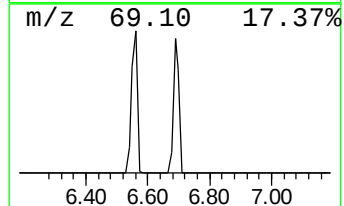
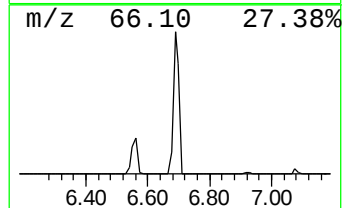
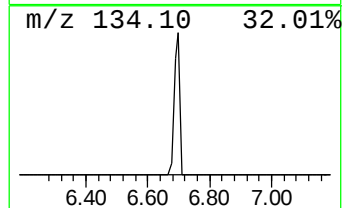
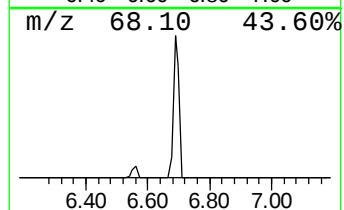
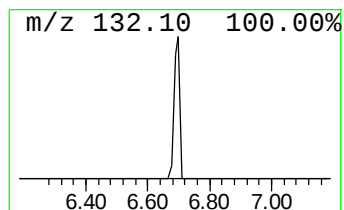
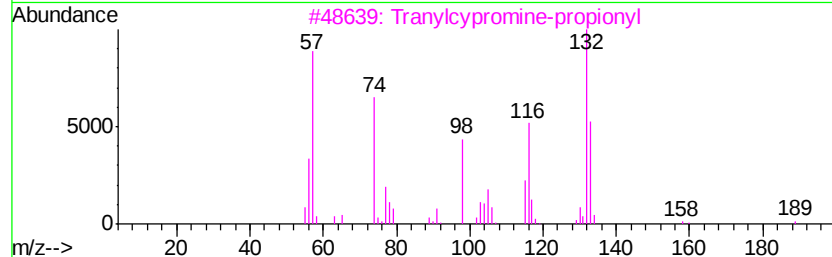
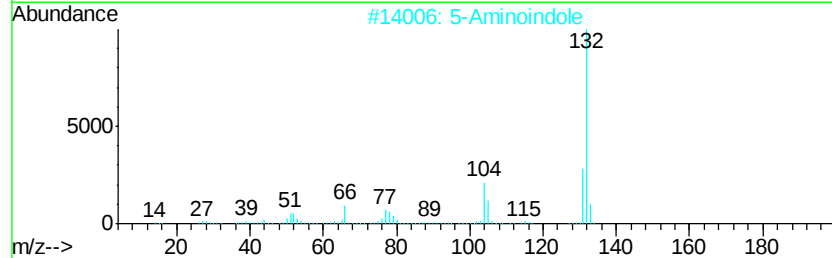
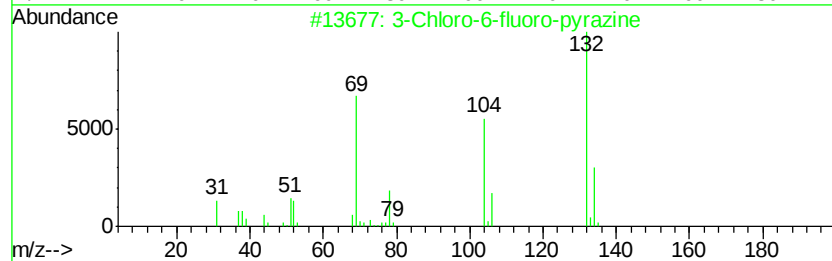
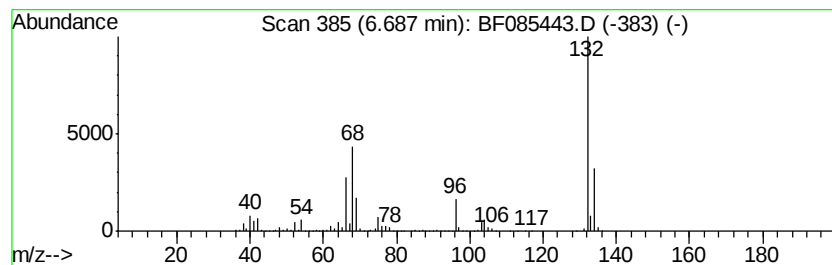
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 unknown6.69 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.69 | 82.78 ng | 4390350 | 1,4-Dichlorobenzene-d4 | 6.93 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 12   |
| 3    |    | Tranylcypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 4    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 9    |
| 5    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

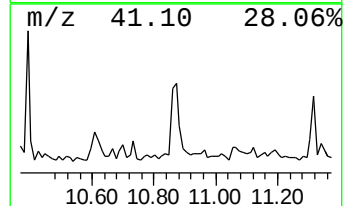
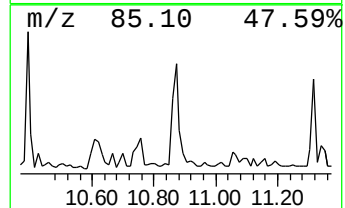
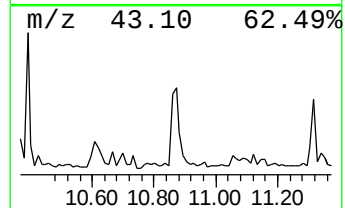
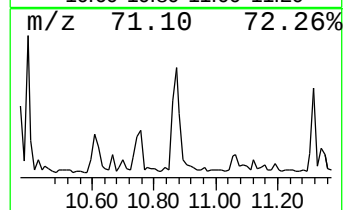
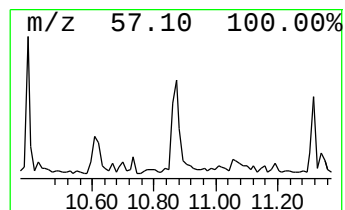
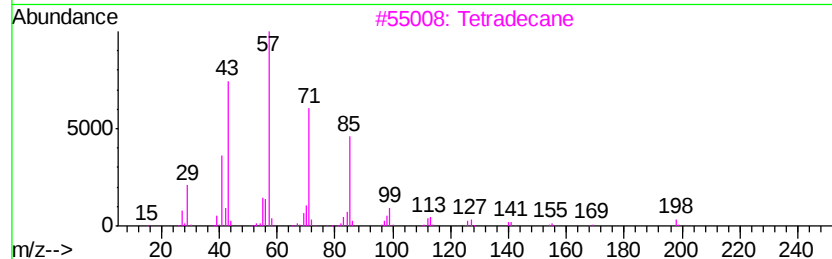
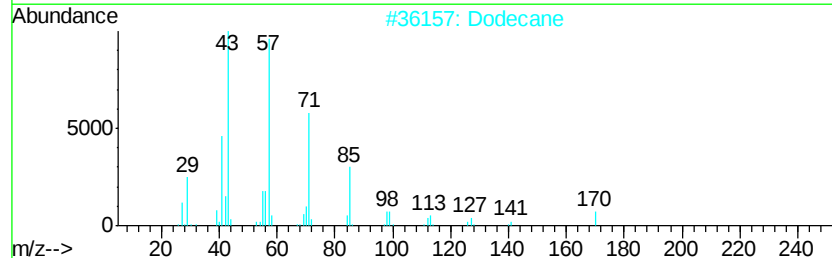
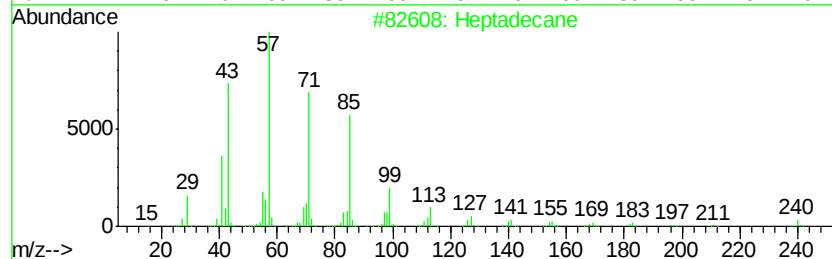
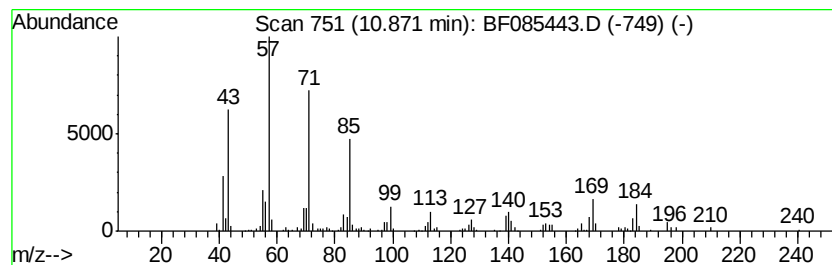
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Heptadecane Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.87 | 3.48 ng | 329243 | Phenanthrene-d10 | 11.45 |

| Hit# of | 5                   | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|---------------------|--------------|-----|---------|-------------|------|
| 1       | Heptadecane         |              | 240 | C17H36  | 000629-78-7 | 95   |
| 2       | Dodecane            |              | 170 | C12H26  | 000112-40-3 | 90   |
| 3       | Tetradecane         |              | 198 | C14H30  | 000629-59-4 | 76   |
| 4       | Tridecane           |              | 184 | C13H28  | 000629-50-5 | 74   |
| 5       | Dodecane, 3-methyl- |              | 184 | C13H28  | 017312-57-1 | 74   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

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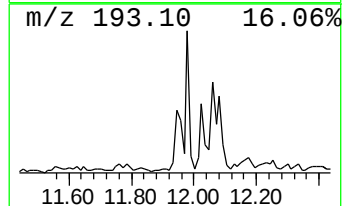
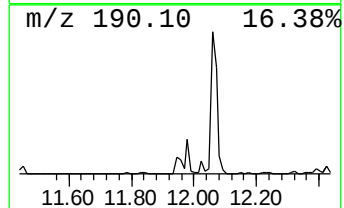
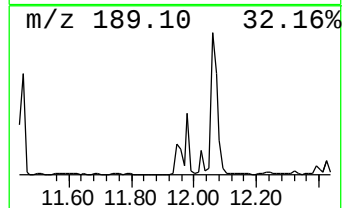
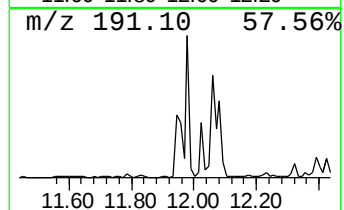
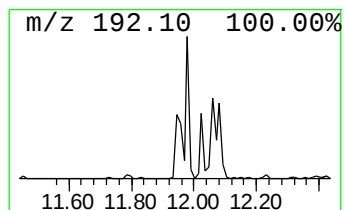
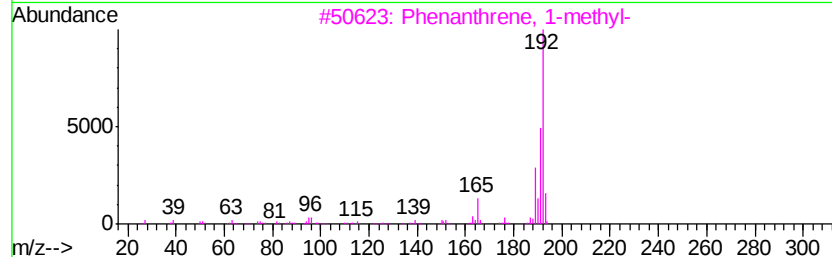
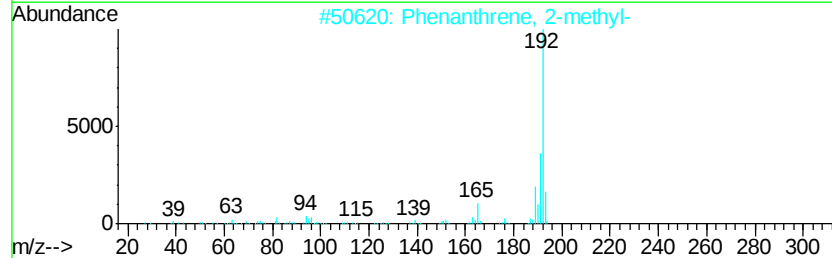
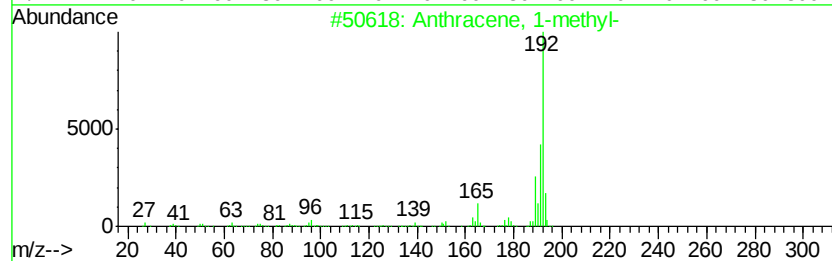
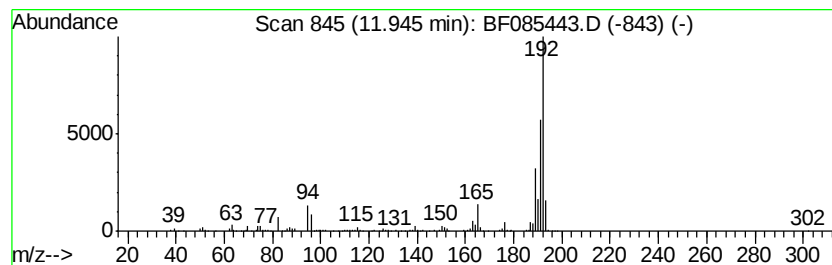
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 Anthracene, 1-methyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.94 | 3.56 ng | 337038 | Phenanthrene-d10 | 11.45 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 95   |
| 2    |    | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 94   |
| 3    |    | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 91   |
| 4    |    | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |
| 5    |    | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

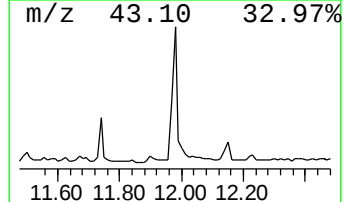
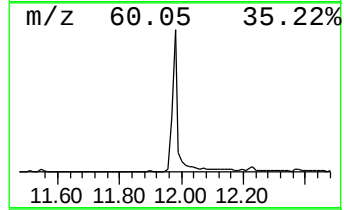
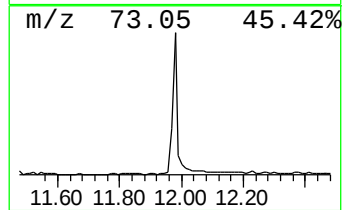
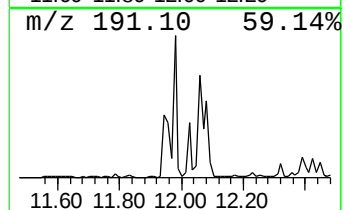
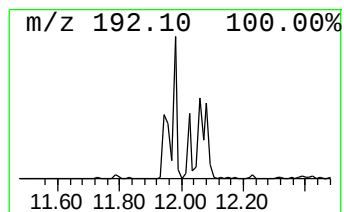
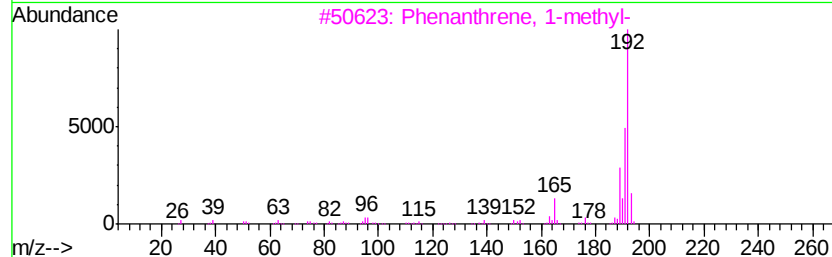
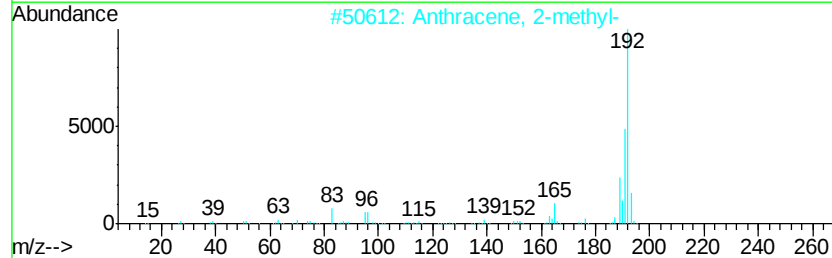
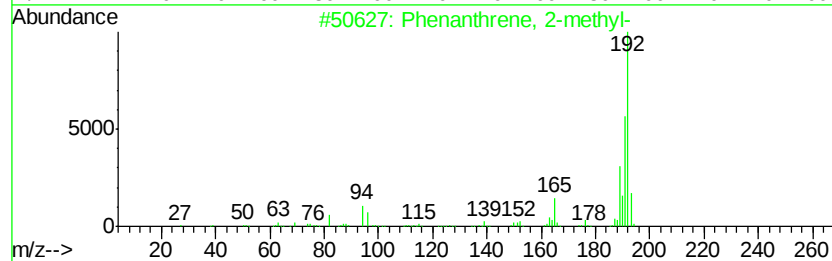
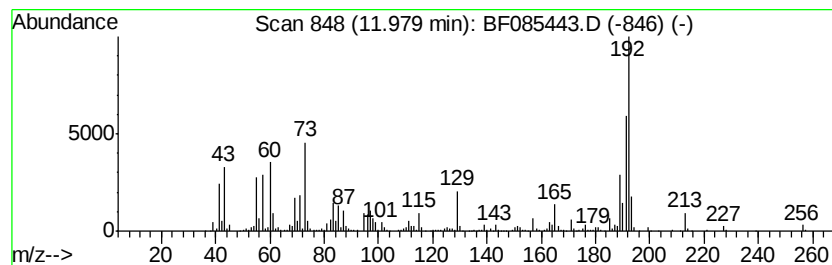
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ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 4

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.98     | 9.60 ng                             | 908126 | Phenanthrene-d10 | 11.45       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 98   |
| 2         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 96   |
| 3         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 96   |
| 4         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 96   |
| 5         | Anthracene, 9-methyl-               | 192    | C15H12           | 000779-02-2 | 92   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

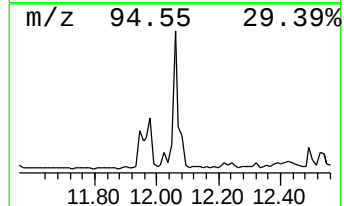
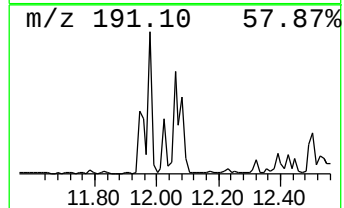
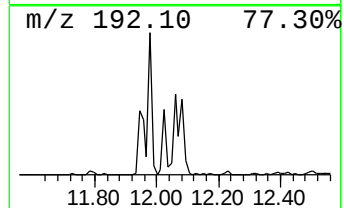
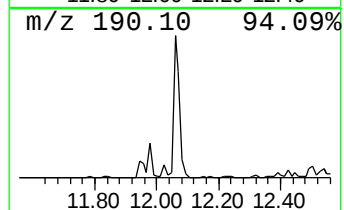
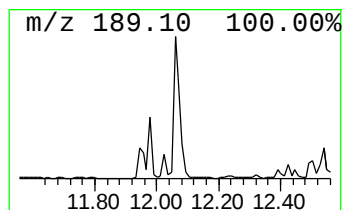
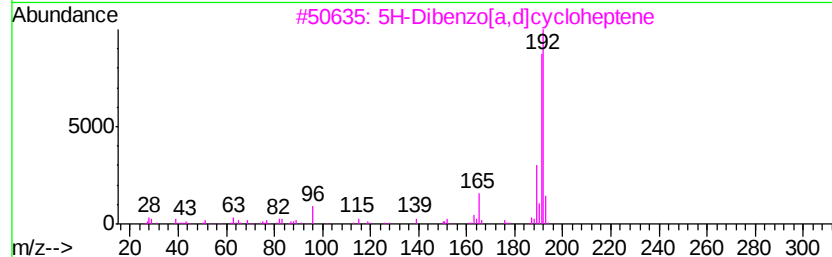
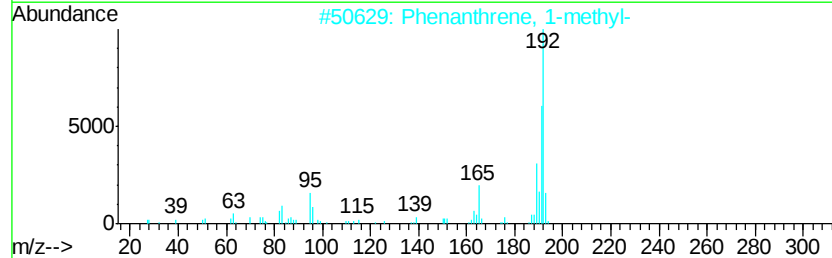
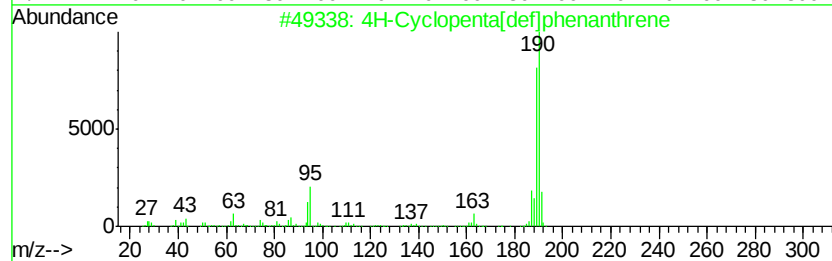
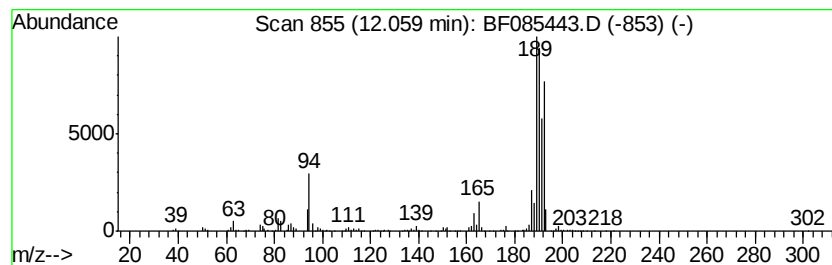
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ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

| R.T.      | EstConc                        | Area    | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------|---------|------------------|-------------|------|
| 12.06     | 10.67 ng                       | 1009000 | Phenanthrene-d10 | 11.45       |      |
| Hit# of 5 | Tentative ID                   | MW      | MolForm          | CAS#        | Qual |
| 1         | 4H-Cyclopenta[def]phenanthrene | 190     | C15H10           | 000203-64-5 | 50   |
| 2         | Phenanthrene, 1-methyl-        | 192     | C15H12           | 000832-69-9 | 42   |
| 3         | 5H-Dibenzo[a,d]cycloheptene    | 192     | C15H12           | 000256-81-5 | 42   |
| 4         | Phenanthrene, 4-methyl-        | 192     | C15H12           | 000832-64-4 | 38   |
| 5         | 1H-Indene, 2-phenyl-           | 192     | C15H12           | 004505-48-0 | 38   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

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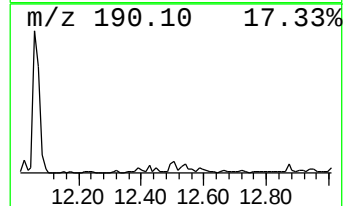
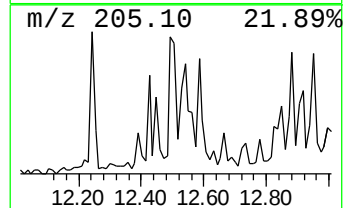
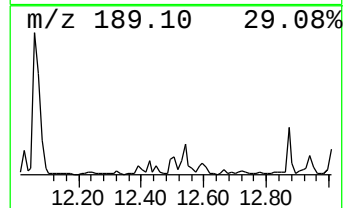
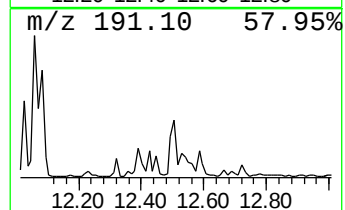
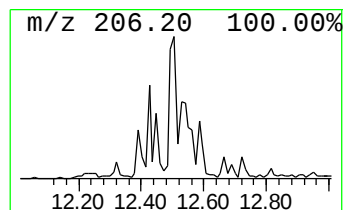
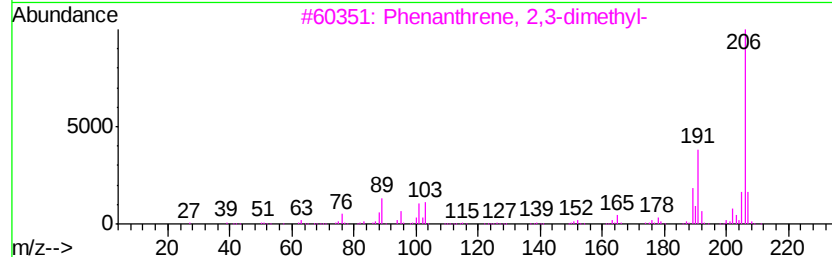
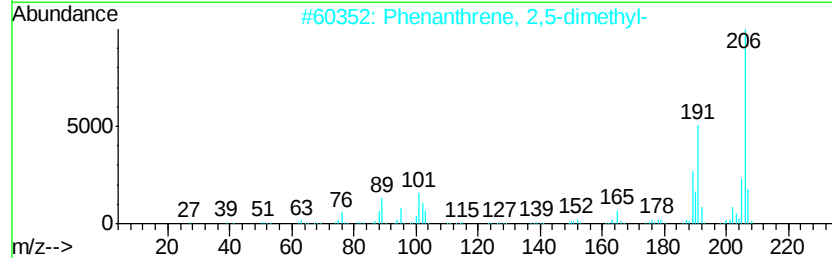
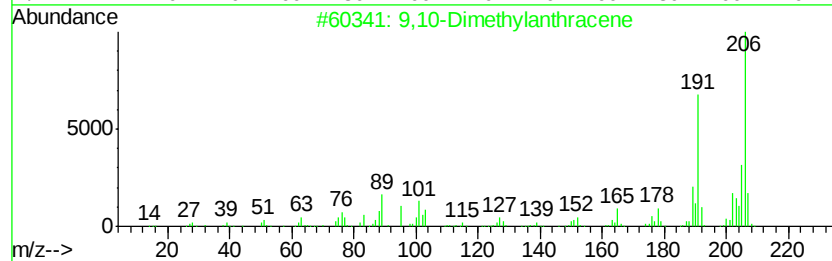
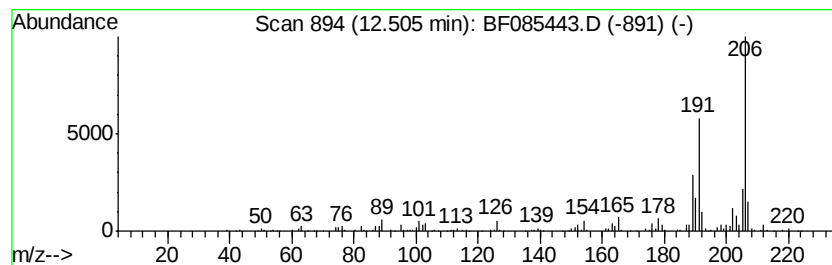
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 9,10-Dimethylantracene Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.51 | 3.31 ng | 312829 | Phenanthrene-d10 | 11.45 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|---------|---|-----------------------------|-----|---------|-------------|------|
| 1       |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 93   |
| 2       |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 93   |
| 3       |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 93   |
| 4       |   | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 87   |
| 5       |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

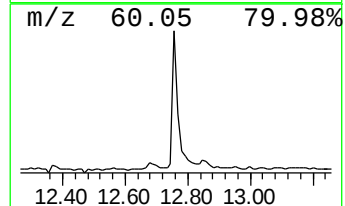
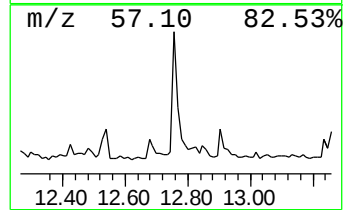
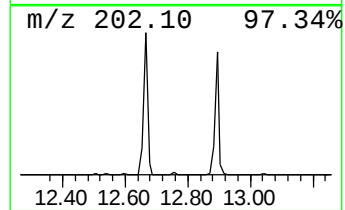
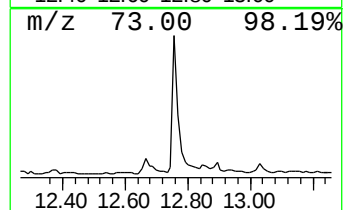
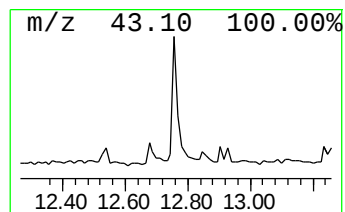
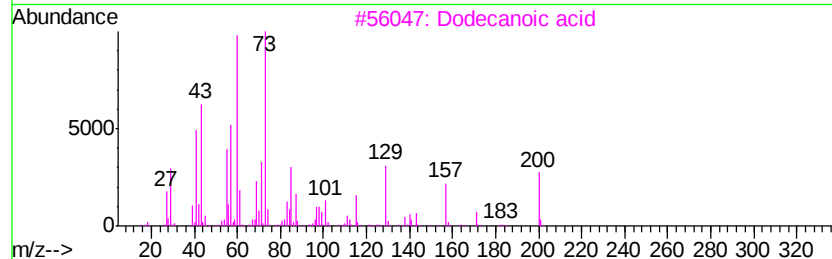
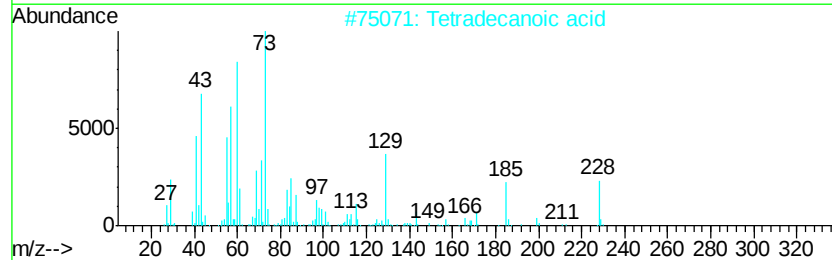
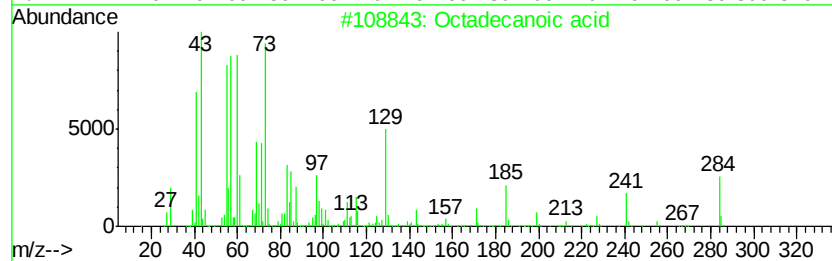
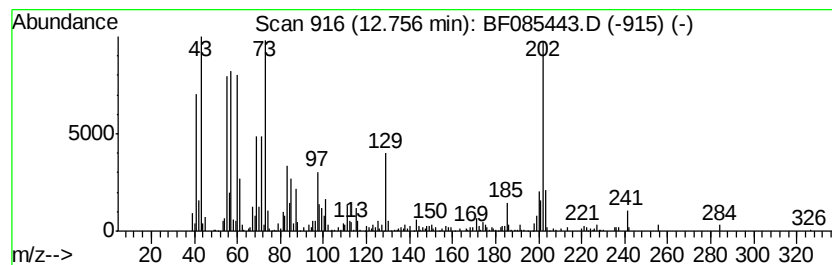
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 Octadecanoic acid Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.76 | 3.31 ng | 313213 | Phenanthrene-d10 | 11.45 |

| Hit# of 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|-----------|--------------------|-----|----------|-------------|------|
| 1         | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2         | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 55   |
| 3         | Dodecanoic acid    | 200 | C12H24O2 | 000143-07-7 | 55   |
| 4         | n-Decanoic acid    | 172 | C10H20O2 | 000334-48-5 | 49   |
| 5         | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 46   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

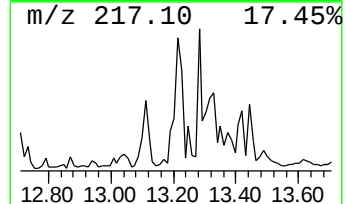
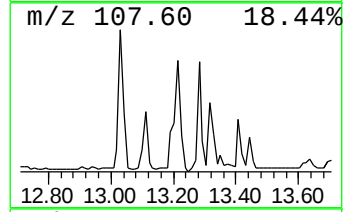
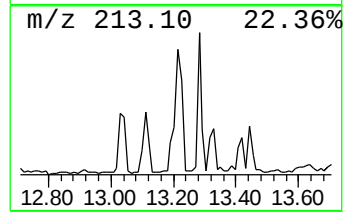
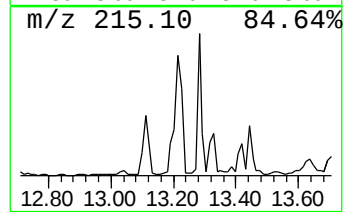
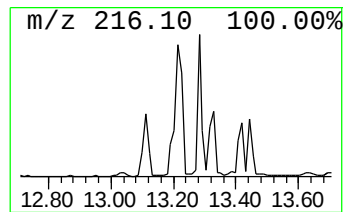
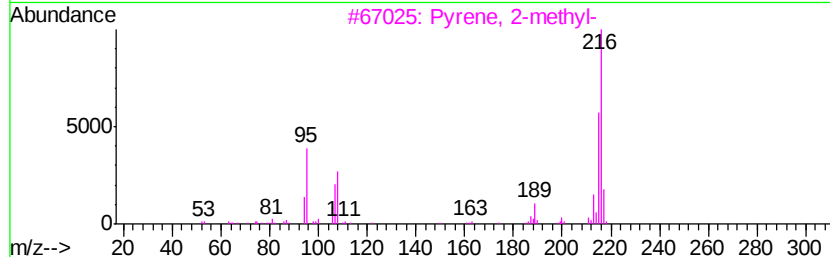
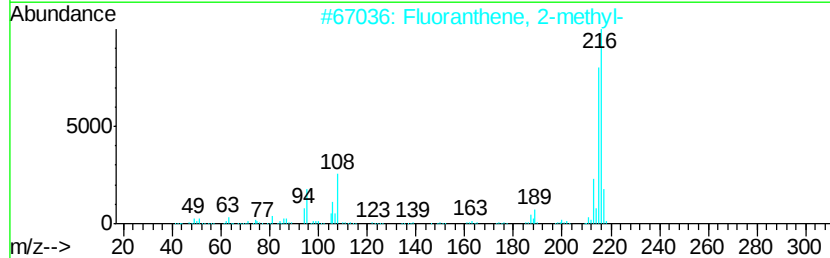
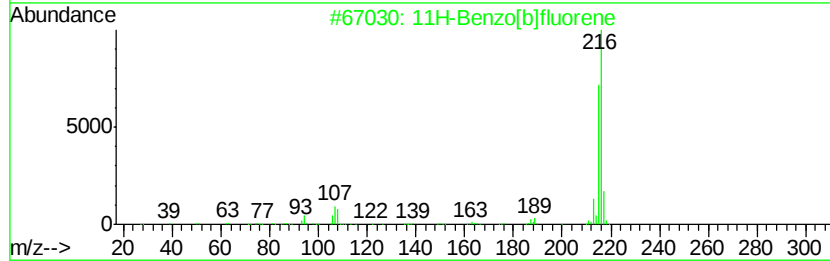
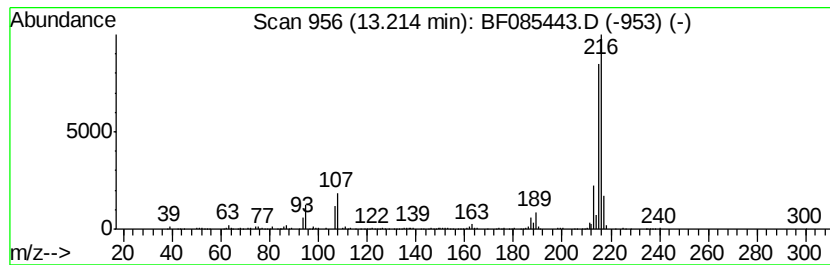
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.21 | 4.52 ng | 544595 | Chrysene-d12     | 14.09 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 95   |
| 2    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 95   |
| 3    |    | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 95   |
| 4    |    | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 93   |
| 5    |    | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 91   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

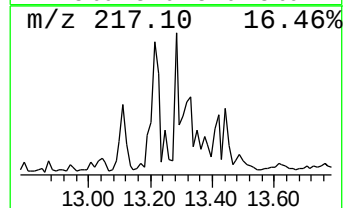
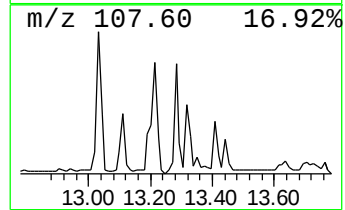
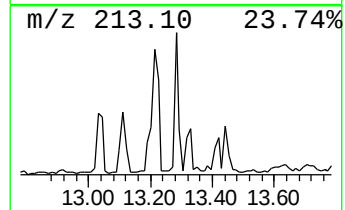
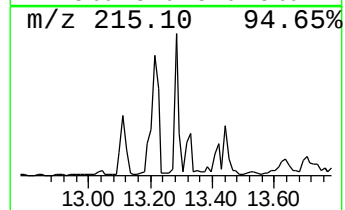
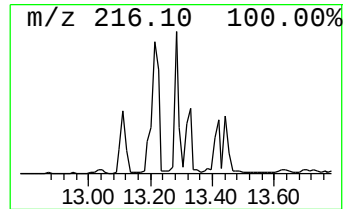
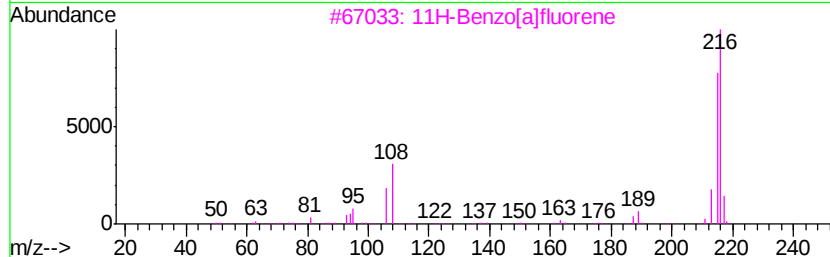
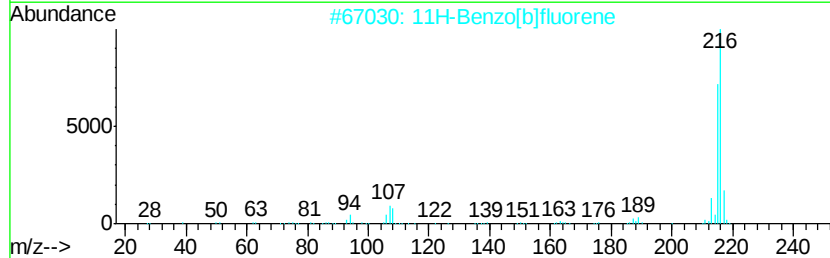
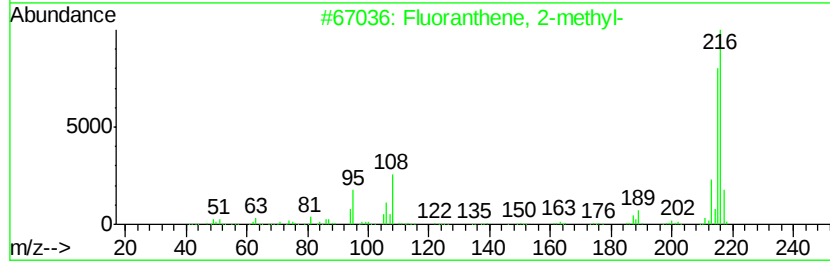
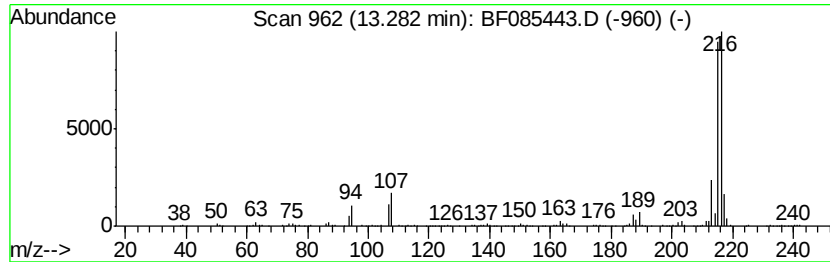
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Fluoranthene, 2-methyl- Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.28 | 2.92 ng | 351136 | Chrysene-d12     | 14.09 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 94   |
| 2    |    | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 93   |
| 3    |    | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 91   |
| 4    |    | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 89   |
| 5    |    | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 83   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

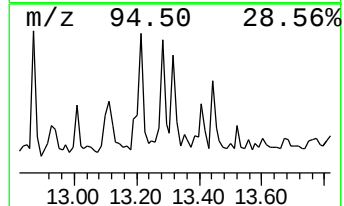
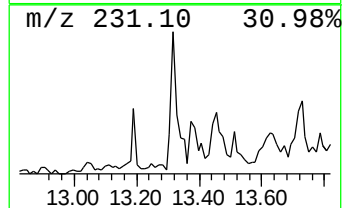
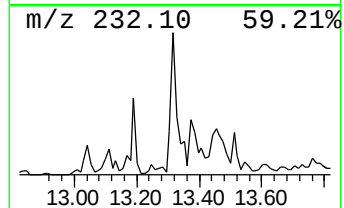
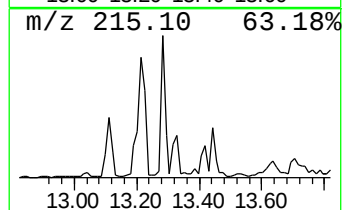
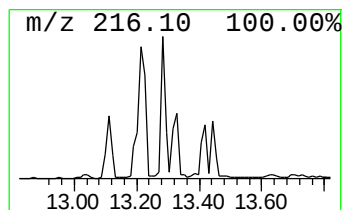
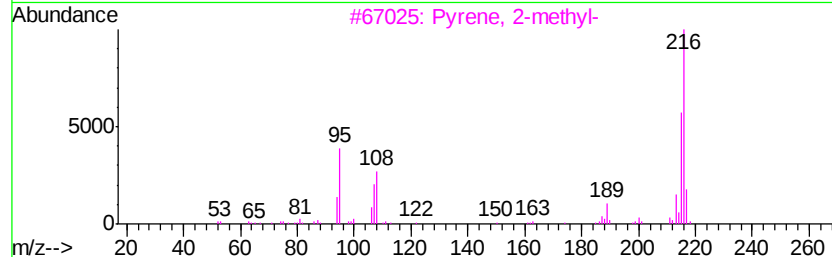
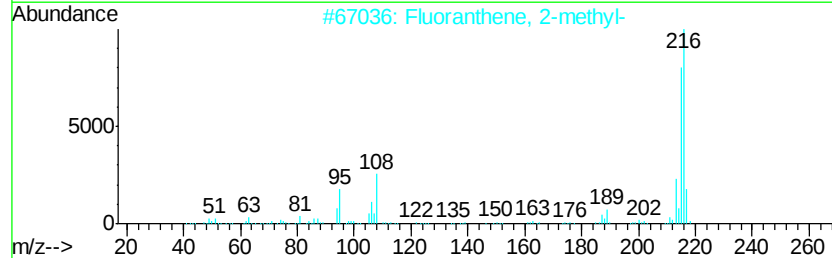
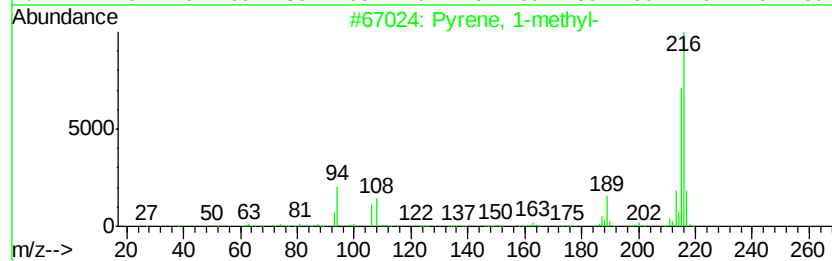
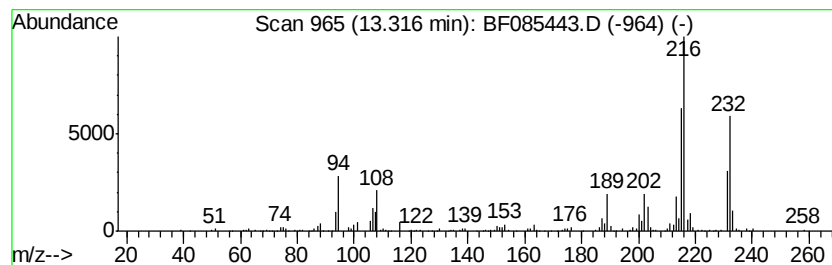
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Pyrene, 1-methyl- Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.32 | 2.56 ng | 308376 | Chrysene-d12     | 14.09 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 91   |
| 2    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 55   |
| 3    |    | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 50   |
| 4    |    | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 46   |
| 5    |    | Pyrene, 4-methyl-       | 216 | C17H12  | 003353-12-6 | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

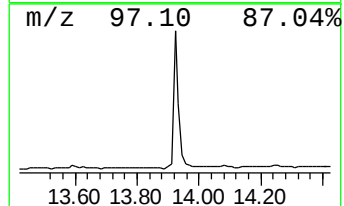
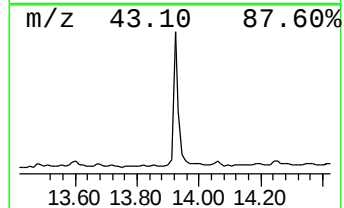
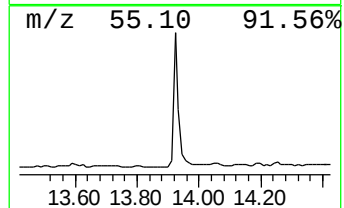
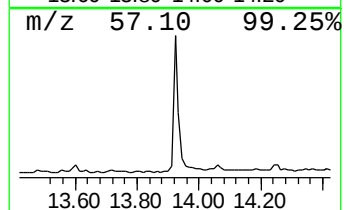
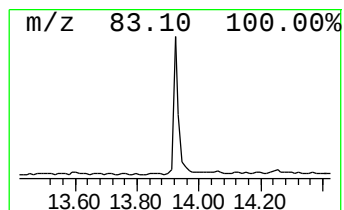
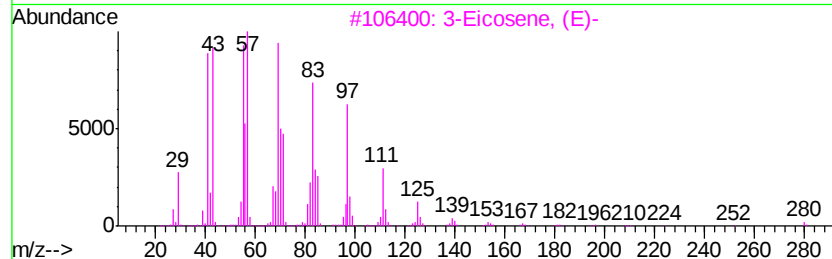
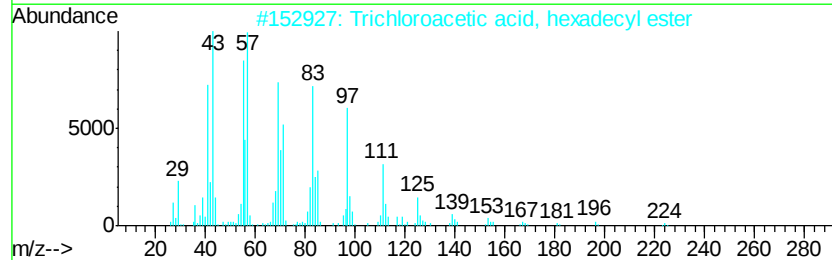
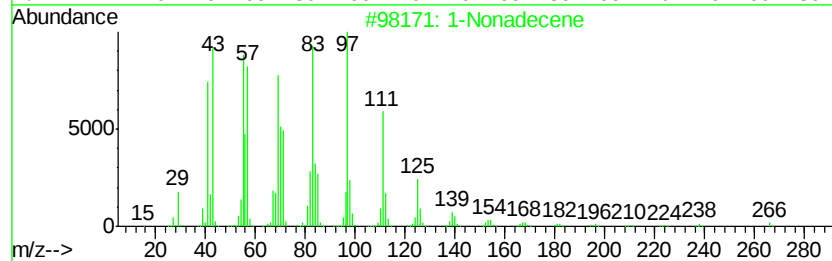
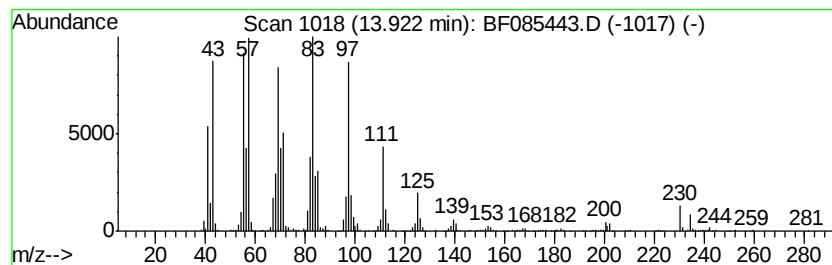
Instrument :  
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ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 1-Nonadecene Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 13.92     | 6.96 ng                             | 838141 | Chrysene-d12     | 14.09       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 1-Nonadecene                        | 266    | C19H38           | 018435-45-5 | 90   |
| 2         | Trichloroacetic acid, hexadecyl ... | 386    | C18H33Cl3O2      | 074339-54-1 | 90   |
| 3         | 3-Eicosene, (E)-                    | 280    | C20H40           | 074685-33-9 | 90   |
| 4         | E-14-Hexadecenal                    | 238    | C16H30O          | 330207-53-9 | 89   |
| 5         | Trichloroacetic acid, pentadecyl... | 372    | C17H31Cl3O2      | 074339-53-0 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

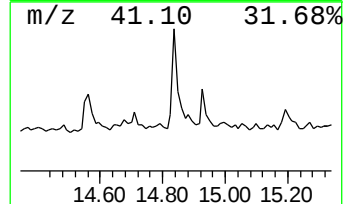
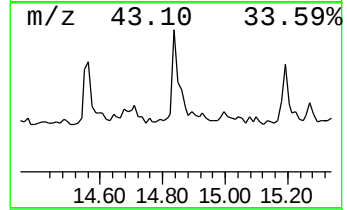
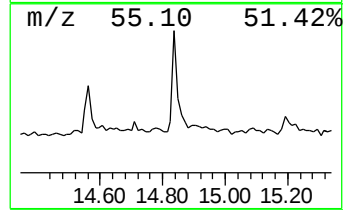
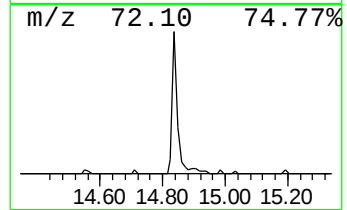
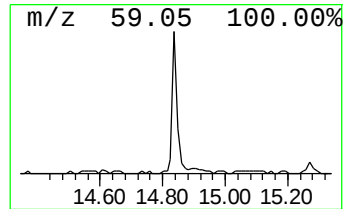
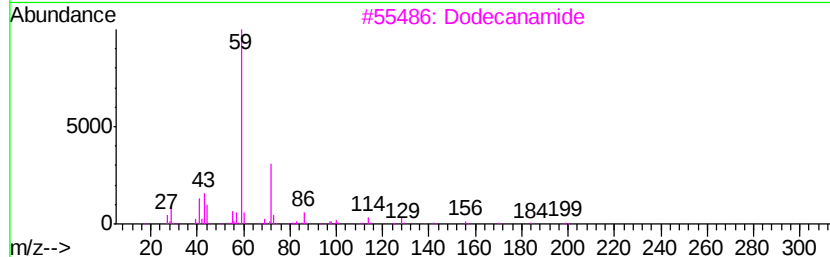
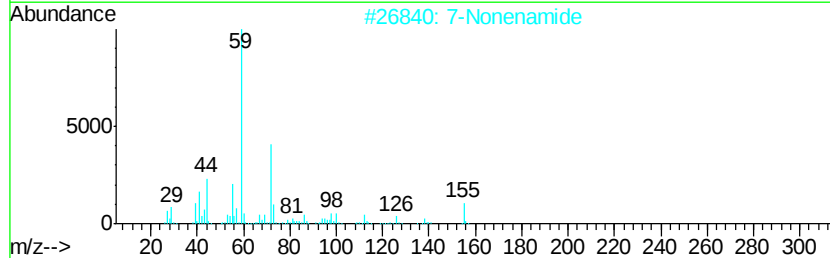
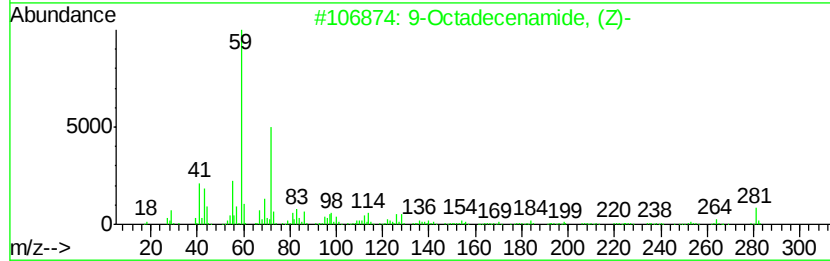
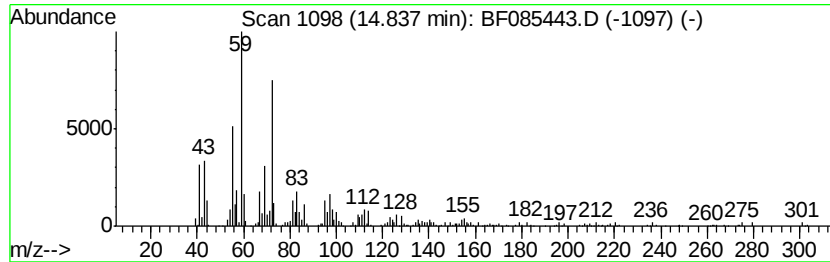
Instrument :  
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ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 9-Octadecenamide, (Z)- Concentration Rank 19

| R.T.      | EstConc                | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------|--------|------------------|-------------|------|
| 14.84     | 2.66 ng                | 155966 | Perylene-d12     | 15.55       |      |
| Hit# of 5 | Tentative ID           | MW     | MolForm          | CAS#        | Qual |
| 1         | 9-Octadecenamide, (Z)- | 281    | C18H35NO         | 000301-02-0 | 86   |
| 2         | 7-Nonenamide           | 155    | C9H17NO          | 090949-53-4 | 58   |
| 3         | Dodecanamide           | 199    | C12H25NO         | 001120-16-7 | 43   |
| 4         | Nonanamide             | 157    | C9H19NO          | 001120-07-6 | 43   |
| 5         | Pentanamide, 4-methyl- | 115    | C6H13NO          | 001119-29-5 | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

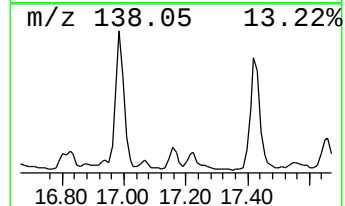
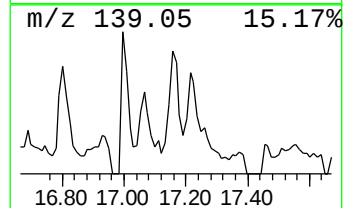
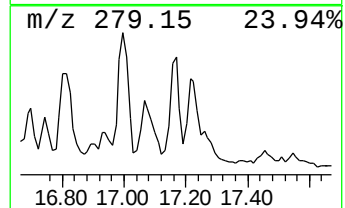
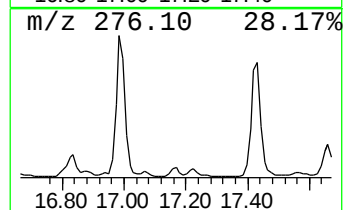
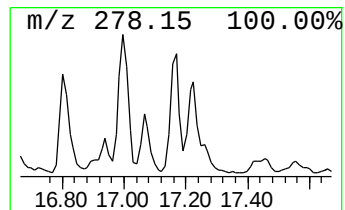
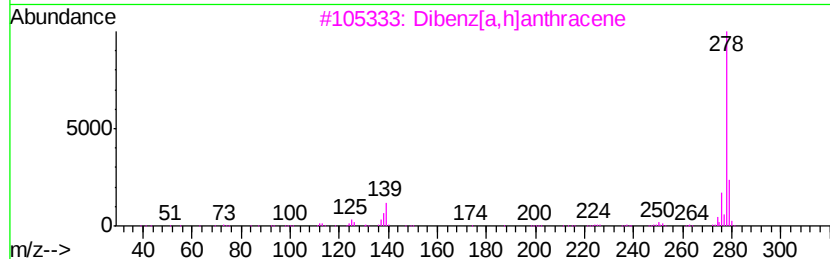
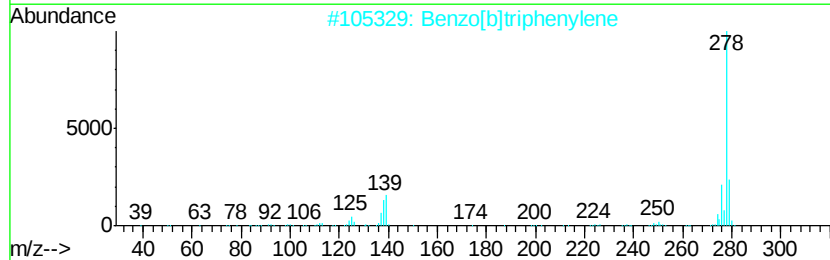
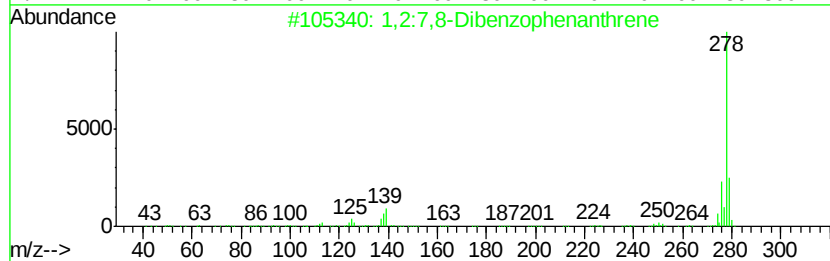
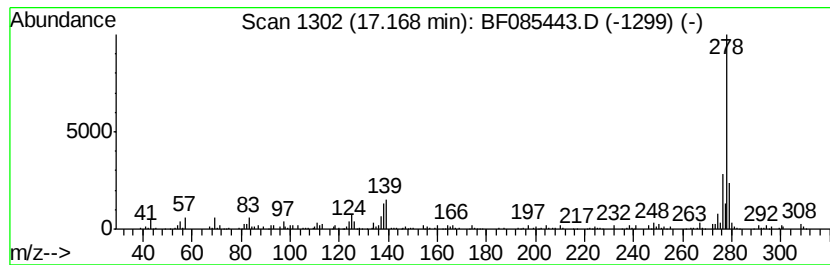
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 1,2:7,8-Dibenzophenanthrene Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.17 | 2.68 ng | 156924 | Perylene-d12     | 15.55 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | 1,2:7,8-Dibenzophenanthrene | 278 | C22H14  | 000213-46-7 | 97   |
| 2    |    | Benzo[b]triphenylene        | 278 | C22H14  | 000215-58-7 | 96   |
| 3    |    | Dibenz[a,h]anthracene       | 278 | C22H14  | 000053-70-3 | 96   |
| 4    |    | Benzo[b]chrysene            | 278 | C22H14  | 000214-17-5 | 95   |
| 5    |    | Pentacene                   | 278 | C22H14  | 000135-48-8 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

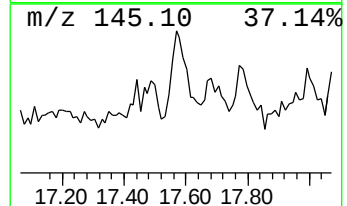
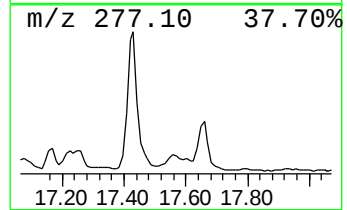
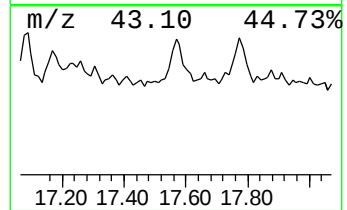
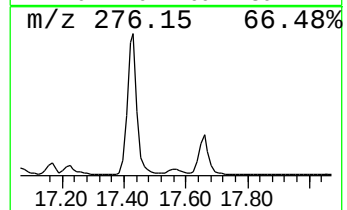
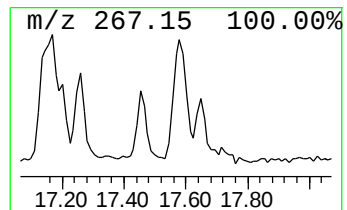
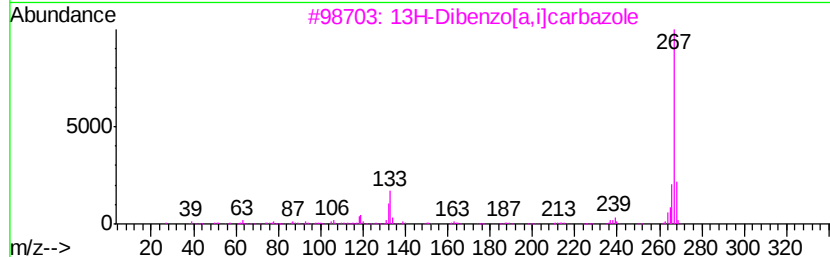
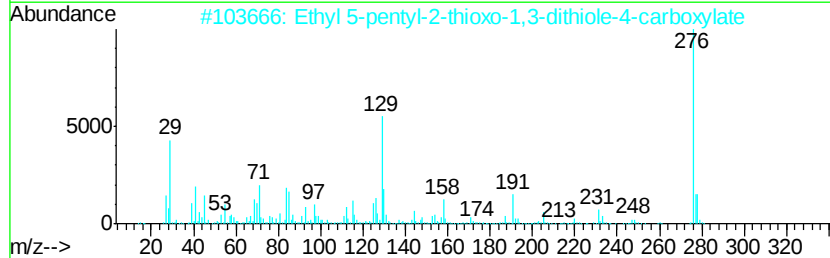
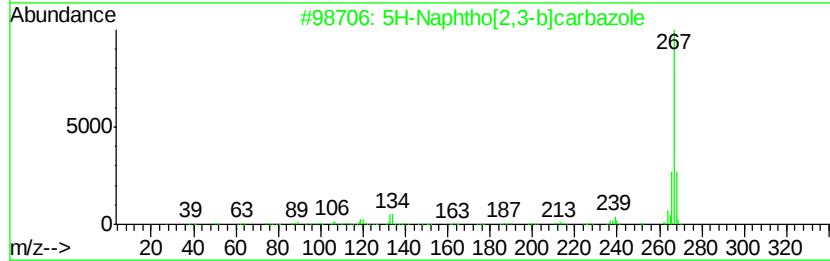
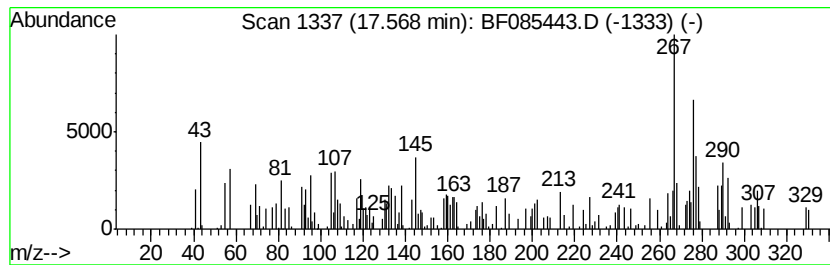
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown17.57 Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.57 | 2.99 ng | 175146 | Perylene-d12     | 15.55 |

| Hit# | of                                  | 5       | Tentative ID | MW  | MolForm     | CAS#        | Qual |
|------|-------------------------------------|---------|--------------|-----|-------------|-------------|------|
| 1    | 5H-Naphtho[2,3-b]carbazole          |         |              | 267 | C20H13N     | 000248-96-4 | 25   |
| 2    | Ethyl 5-pentyl-2-thioxo-1,3-dith... |         |              | 276 | C11H16O2S3  | 120356-16-3 | 15   |
| 3    | 13H-Dibenzo[a,i]carbazole           |         |              | 267 | C20H13N     | 000239-64-5 | 15   |
| 4    | 2-Chloro-7-methoxyphenazine         | 5,10... |              | 276 | C13H9ClN2O3 | 023677-04-5 | 11   |
| 5    | 5H-Naphtho[2,3-c]carbazole          |         |              | 267 | C20H13N     | 000198-96-9 | 11   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
Data File : BF085443.D  
Acq On : 14 Mar 2016 21:53  
Operator : UM/SJ  
Sample : H1722-10  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

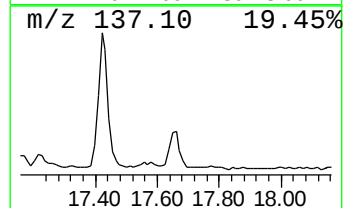
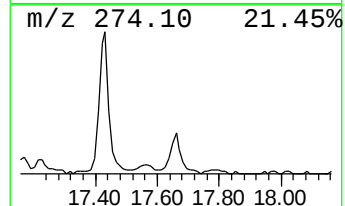
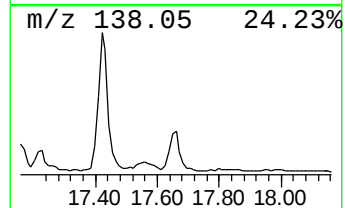
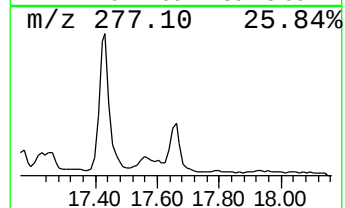
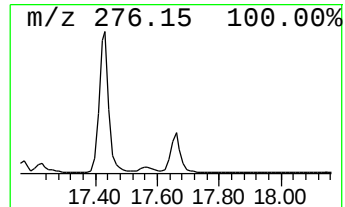
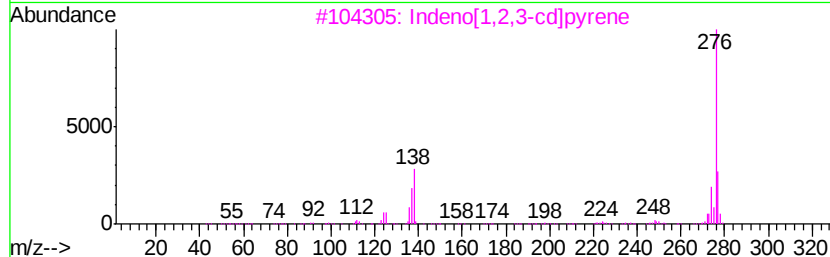
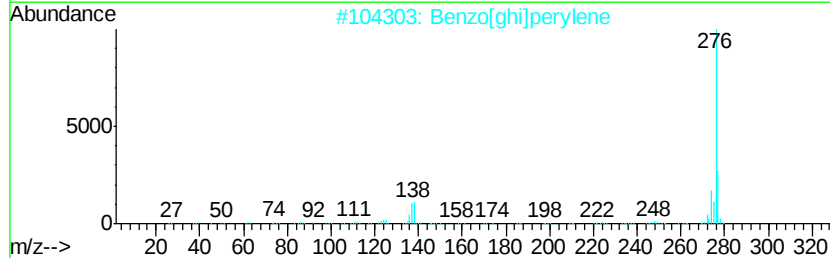
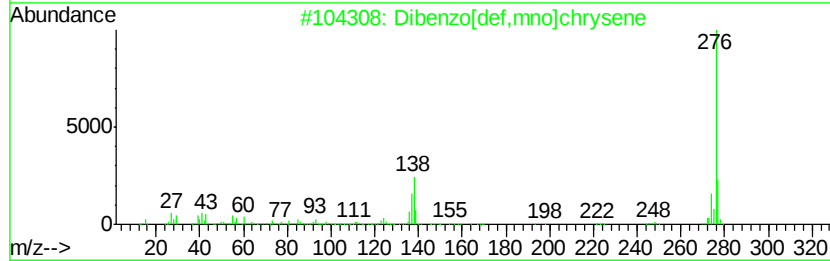
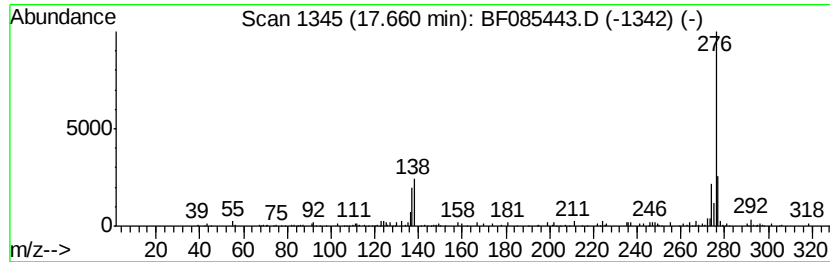
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Dibenzo[def,mno]chrysene Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.66 | 2.84 ng | 166401 | Perylene-d12     | 15.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Dibenzo[def,mno]chrysene            | 276 | C22H12   | 000191-26-4 | 98   |
| 2    |    | Benzo[ghi]perylene                  | 276 | C22H12   | 000191-24-2 | 93   |
| 3    |    | Indeno[1,2,3-cd]pyrene              | 276 | C22H12   | 000193-39-5 | 91   |
| 4    |    | Indeno[1,2,3-cd]fluoranthene        | 276 | C22H12   | 000193-43-1 | 81   |
| 5    |    | 1H-Cyclopenta[4,5]pyrido[1,2-a]b... | 276 | C17H16N4 | 329707-31-5 | 47   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF031416\  
 Data File : BF085443.D  
 Acq On : 14 Mar 2016 21:53  
 Operator : UM/SJ  
 Sample : H1722-10  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 W170TH-SB-22(0.5-2.0)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF030316.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT 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-Pentanone, 4-hy...  | 5.13  | 5.8 ng  |       | 306428   | 1                     | 6.93  | 1060750 | 20.0 |
| unknown6.69           | 6.69  | 82.8 ng |       | 4390350  | 1                     | 6.93  | 1060750 | 20.0 |
| Heptadecane           | 10.87 | 3.5 ng  |       | 329243   | 4                     | 11.45 | 1891280 | 20.0 |
| Anthracene, 1-met...  | 11.94 | 3.6 ng  |       | 337038   | 4                     | 11.45 | 1891280 | 20.0 |
| Phenanthrene, 2-m...  | 11.98 | 9.6 ng  |       | 908126   | 4                     | 11.45 | 1891280 | 20.0 |
| 4H-Cyclopenta[def...  | 12.06 | 10.7 ng |       | 1009000  | 4                     | 11.45 | 1891280 | 20.0 |
| 9,10-Dimethylanth...  | 12.51 | 3.3 ng  |       | 312829   | 4                     | 11.45 | 1891280 | 20.0 |
| Octadecanoic acid     | 12.76 | 3.3 ng  |       | 313213   | 4                     | 11.45 | 1891280 | 20.0 |
| 11H-Benzo[b]fluorene  | 13.21 | 4.5 ng  |       | 544595   | 5                     | 14.09 | 2408510 | 20.0 |
| Fluoranthene, 2-m...  | 13.28 | 2.9 ng  |       | 351136   | 5                     | 14.09 | 2408510 | 20.0 |
| Pyrene, 1-methyl-     | 13.32 | 2.6 ng  |       | 308376   | 5                     | 14.09 | 2408510 | 20.0 |
| 1-Nonadecene          | 13.92 | 7.0 ng  |       | 838141   | 5                     | 14.09 | 2408510 | 20.0 |
| 9-Octadecenamide, ... | 14.84 | 2.7 ng  |       | 155966   | 6                     | 15.55 | 1172010 | 20.0 |
| 1,2:7,8-Dibenzoph...  | 17.17 | 2.7 ng  |       | 156924   | 6                     | 15.55 | 1172010 | 20.0 |
| unknown17.57          | 17.57 | 3.0 ng  |       | 175146   | 6                     | 15.55 | 1172010 | 20.0 |
| Dibenzo[def,mno]c...  | 17.66 | 2.8 ng  |       | 166401   | 6                     | 15.55 | 1172010 | 20.0 |