

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
 Data File : BF089978.D  
 Acq On : 2 Sep 2016 16:43  
 Operator : UM/SJ  
 Sample : H4683-01  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SOILSAMPLE-1

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-BF083116.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         | 1.644       | 4             | 5           | 13           | rVB      | 52678          | 100635        | 2.76%           | 0.146%        |
| 2         | 1.758       | 13            | 15          | 55           | rVB      | 1262789        | 3491668       | 95.88%          | 5.057%        |
| 3         | 4.753       | 274           | 277         | 283          | rBV      | 549670         | 673286        | 18.49%          | 0.975%        |
| 4         | 5.210       | 314           | 317         | 319          | rBV      | 2276163        | 2714823       | 74.55%          | 3.932%        |
| 5         | 6.273       | 407           | 410         | 413          | rBV      | 2826176        | 2790151       | 76.62%          | 4.041%        |
| 6         | 6.399       | 418           | 421         | 423          | rBV      | 3258612        | 2936373       | 80.63%          | 4.253%        |
| 7         | 6.639       | 439           | 442         | 444          | rBV      | 1092145        | 1166706       | 32.04%          | 1.690%        |
| 8         | 6.787       | 453           | 455         | 458          | rVB      | 2075118        | 2069819       | 56.84%          | 2.998%        |
| 9         | 7.199       | 488           | 491         | 494          | rBV      | 1991970        | 1775343       | 48.75%          | 2.571%        |
| 10        | 7.919       | 552           | 554         | 558          | rBV      | 1751703        | 1506369       | 41.36%          | 2.182%        |
| 11        | 8.730       | 622           | 625         | 627          | rBV      | 88714          | 81715         | 2.24%           | 0.118%        |
| 12        | 8.993       | 646           | 648         | 651          | rBV      | 2904198        | 3604269       | 98.97%          | 5.220%        |
| 13        | 9.325       | 675           | 677         | 678          | rBV      | 141883         | 135607        | 3.72%           | 0.196%        |
| 14        | 9.393       | 681           | 683         | 685          | rBV      | 996079         | 792260        | 21.76%          | 1.147%        |
| 15        | 9.439       | 685           | 687         | 690          | rVB      | 68497          | 87346         | 2.40%           | 0.127%        |
| 16        | 9.519       | 692           | 694         | 697          | rVB      | 218981         | 230910        | 6.34%           | 0.334%        |
| 17        | 9.668       | 704           | 707         | 708          | rBV      | 1647894        | 1649182       | 45.29%          | 2.389%        |
| 18        | 9.690       | 708           | 709         | 712          | rVB      | 187873         | 191354        | 5.25%           | 0.277%        |
| 19        | 9.862       | 722           | 724         | 726          | rBV      | 121577         | 153747        | 4.22%           | 0.223%        |
| 20        | 9.976       | 732           | 734         | 736          | rBV      | 62321          | 84232         | 2.31%           | 0.122%        |
| 21        | 10.068      | 740           | 742         | 745          | rBV2     | 61791          | 134979        | 3.71%           | 0.195%        |
| 22        | 10.113      | 745           | 746         | 748          | rVB      | 121074         | 103447        | 2.84%           | 0.150%        |
| 23        | 10.159      | 748           | 750         | 752          | rBV3     | 36889          | 86676         | 2.38%           | 0.126%        |
| 24        | 10.205      | 752           | 754         | 756          | rBV      | 243380         | 186091        | 5.11%           | 0.270%        |
| 25        | 10.273      | 756           | 760         | 763          | rBV3     | 78799          | 126267        | 3.47%           | 0.183%        |
| 26        | 10.388      | 769           | 770         | 771          | rVB      | 284564         | 195211        | 5.36%           | 0.283%        |
| 27        | 10.445      | 771           | 775         | 777          | rBV      | 2152983        | 2028055       | 55.69%          | 2.937%        |
| 28        | 10.582      | 782           | 787         | 790          | rBV3     | 123291         | 267726        | 7.35%           | 0.388%        |
| 29        | 10.753      | 798           | 802         | 803          | rBV      | 86225          | 159792        | 4.39%           | 0.231%        |
| 30        | 10.902      | 811           | 815         | 816          | rBV2     | 93008          | 166121        | 4.56%           | 0.241%        |
| 31        | 10.936      | 816           | 818         | 821          | rVB      | 159186         | 200814        | 5.51%           | 0.291%        |
| 32        | 10.993      | 821           | 823         | 824          | rBV      | 86444          | 99537         | 2.73%           | 0.144%        |
| 33        | 11.028      | 824           | 826         | 833          | rVB      | 230074         | 301987        | 8.29%           | 0.437%        |
| 34        | 11.165      | 833           | 838         | 839          | rBV2     | 1715614        | 3641667       | 100.00%         | 5.274%        |

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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-BF083116.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 11.211 | 839  | 842  | 844  | rVB  | 559886  | 543553  | 14.93% | 0.787% |
| 36 | 11.313 | 847  | 851  | 852  | rVB3 | 80158   | 136923  | 3.76%  | 0.198% |
| 37 | 11.359 | 852  | 855  | 857  | rBV2 | 138552  | 233501  | 6.41%  | 0.338% |
| 38 | 11.439 | 860  | 862  | 863  | rBV  | 65360   | 83915   | 2.30%  | 0.122% |
| 39 | 11.462 | 863  | 864  | 866  | rVB  | 203785  | 164538  | 4.52%  | 0.238% |
| 40 | 11.542 | 868  | 871  | 873  | rVB2 | 69794   | 135504  | 3.72%  | 0.196% |
| 41 | 11.633 | 877  | 879  | 880  | rBV  | 661363  | 559012  | 15.35% | 0.810% |
| 42 | 11.656 | 880  | 881  | 883  | rVB  | 487118  | 594699  | 16.33% | 0.861% |
| 43 | 11.713 | 883  | 886  | 887  | rBV  | 158745  | 234555  | 6.44%  | 0.340% |
| 44 | 11.748 | 887  | 889  | 893  | rVB2 | 571203  | 1216465 | 33.40% | 1.762% |
| 45 | 11.839 | 893  | 897  | 898  | rBV3 | 58472   | 129622  | 3.56%  | 0.188% |
| 46 | 11.862 | 898  | 899  | 901  | rVB2 | 98552   | 81260   | 2.23%  | 0.118% |
| 47 | 11.931 | 901  | 905  | 910  | rVB2 | 375439  | 664984  | 18.26% | 0.963% |
| 48 | 12.079 | 915  | 918  | 919  | rBV2 | 189959  | 237750  | 6.53%  | 0.344% |
| 49 | 12.102 | 919  | 920  | 922  | rBV  | 71630   | 104581  | 2.87%  | 0.151% |
| 50 | 12.182 | 924  | 927  | 928  | rBV  | 432708  | 461622  | 12.68% | 0.669% |
| 51 | 12.216 | 928  | 930  | 931  | rVV  | 203452  | 213170  | 5.85%  | 0.309% |
| 52 | 12.262 | 931  | 934  | 938  | rVB3 | 322181  | 482755  | 13.26% | 0.699% |
| 53 | 12.342 | 938  | 941  | 943  | rBV  | 2558647 | 2790279 | 76.62% | 4.041% |
| 54 | 12.399 | 943  | 946  | 947  | rBV2 | 89461   | 134742  | 3.70%  | 0.195% |
| 55 | 12.422 | 947  | 948  | 950  | rVB  | 127641  | 163796  | 4.50%  | 0.237% |
| 56 | 12.468 | 950  | 952  | 956  | rBV3 | 172137  | 349270  | 9.59%  | 0.506% |
| 57 | 12.559 | 958  | 960  | 963  | rBV  | 2270620 | 2782904 | 76.42% | 4.031% |
| 58 | 12.616 | 963  | 965  | 968  | rVB2 | 144830  | 261265  | 7.17%  | 0.378% |
| 59 | 12.685 | 968  | 971  | 972  | rBV  | 168216  | 207431  | 5.70%  | 0.300% |
| 60 | 12.719 | 972  | 974  | 976  | rBV  | 3298190 | 3095670 | 85.01% | 4.484% |
| 61 | 12.788 | 976  | 980  | 981  | rBV2 | 235039  | 298049  | 8.18%  | 0.432% |
| 62 | 12.891 | 983  | 989  | 991  | rVB2 | 379723  | 766495  | 21.05% | 1.110% |
| 63 | 12.959 | 991  | 995  | 996  | rBV  | 211742  | 312395  | 8.58%  | 0.452% |
| 64 | 12.994 | 996  | 998  | 1000 | rVV  | 427367  | 442722  | 12.16% | 0.641% |
| 65 | 13.085 | 1004 | 1006 | 1007 | rBV  | 221149  | 225144  | 6.18%  | 0.326% |
| 66 | 13.108 | 1007 | 1008 | 1011 | rBV  | 211788  | 218752  | 6.01%  | 0.317% |
| 67 | 13.188 | 1013 | 1015 | 1020 | rVB3 | 93397   | 201387  | 5.53%  | 0.292% |
| 68 | 13.314 | 1020 | 1026 | 1027 | rBV5 | 119125  | 381237  | 10.47% | 0.552% |
| 69 | 13.394 | 1030 | 1033 | 1035 | rBV  | 242469  | 364024  | 10.00% | 0.527% |
| 70 | 13.496 | 1040 | 1042 | 1044 | rBV  | 320187  | 446981  | 12.27% | 0.647% |
| 71 | 13.531 | 1044 | 1045 | 1046 | rVV  | 227120  | 192040  | 5.27%  | 0.278% |

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Misc :  
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Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Integration Parameters: rteint.p  
Integrator: RTE  
Smoothing : ON Filtering: 5  
Sampling : 1 Min Area: 3 % of largest Peak  
Start Thrs: 0.2 Max Peaks: 100  
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 72  | 13.599 | 1049 | 1051 | 1052 | rVB  | 231889  | 244407  | 6.71%  | 0.354% |
| 73  | 13.634 | 1052 | 1054 | 1061 | rBV4 | 259825  | 342782  | 9.41%  | 0.496% |
| 74  | 13.748 | 1061 | 1064 | 1066 | rBV2 | 1986020 | 3430965 | 94.21% | 4.969% |
| 75  | 13.851 | 1070 | 1073 | 1074 | rBV2 | 126289  | 147710  | 4.06%  | 0.214% |
| 76  | 13.885 | 1074 | 1076 | 1077 | rBV2 | 98420   | 109657  | 3.01%  | 0.159% |
| 77  | 13.977 | 1082 | 1084 | 1085 | rVB2 | 98439   | 118190  | 3.25%  | 0.171% |
| 78  | 14.068 | 1090 | 1092 | 1094 | rBV3 | 75796   | 118552  | 3.26%  | 0.172% |
| 79  | 14.137 | 1094 | 1098 | 1099 | rBV  | 393032  | 492759  | 13.53% | 0.714% |
| 80  | 14.171 | 1099 | 1101 | 1103 | rVV  | 197955  | 174407  | 4.79%  | 0.253% |
| 81  | 14.228 | 1103 | 1106 | 1107 | rVB2 | 149355  | 171681  | 4.71%  | 0.249% |
| 82  | 14.262 | 1107 | 1109 | 1113 | rBV5 | 162635  | 397496  | 10.92% | 0.576% |
| 83  | 14.331 | 1113 | 1115 | 1118 | rVB2 | 95380   | 127314  | 3.50%  | 0.184% |
| 84  | 14.434 | 1122 | 1124 | 1125 | rBV  | 87822   | 104939  | 2.88%  | 0.152% |
| 85  | 14.582 | 1135 | 1137 | 1138 | rBV  | 117332  | 132483  | 3.64%  | 0.192% |
| 86  | 14.662 | 1142 | 1144 | 1148 | rVB3 | 100276  | 199969  | 5.49%  | 0.290% |
| 87  | 14.742 | 1148 | 1151 | 1155 | rBV  | 1166486 | 2196790 | 60.32% | 3.182% |
| 88  | 14.834 | 1156 | 1159 | 1161 | rVV2 | 220453  | 292420  | 8.03%  | 0.424% |
| 89  | 14.994 | 1171 | 1173 | 1175 | rVV  | 715391  | 749135  | 20.57% | 1.085% |
| 90  | 15.051 | 1175 | 1178 | 1180 | rVB  | 882704  | 1121824 | 30.81% | 1.625% |
| 91  | 15.097 | 1180 | 1182 | 1190 | rBV2 | 872501  | 1548514 | 42.52% | 2.243% |
| 92  | 15.337 | 1201 | 1203 | 1204 | rBV  | 76066   | 103198  | 2.83%  | 0.149% |
| 93  | 15.531 | 1218 | 1220 | 1222 | rBV  | 157685  | 232164  | 6.38%  | 0.336% |
| 94  | 16.194 | 1274 | 1278 | 1282 | rVB3 | 142484  | 326272  | 8.96%  | 0.473% |
| 95  | 16.331 | 1287 | 1290 | 1294 | rVV2 | 381751  | 697811  | 19.16% | 1.011% |
| 96  | 16.525 | 1305 | 1307 | 1309 | rBV  | 77720   | 136767  | 3.76%  | 0.198% |
| 97  | 16.697 | 1318 | 1322 | 1327 | rVB  | 313581  | 603960  | 16.58% | 0.875% |
| 98  | 18.068 | 1438 | 1442 | 1448 | rVB4 | 93694   | 293079  | 8.05%  | 0.424% |
| 99  | 18.560 | 1480 | 1485 | 1492 | rVB2 | 105647  | 410924  | 11.28% | 0.595% |
| 100 | 18.708 | 1494 | 1498 | 1502 | rBV  | 49023   | 170722  | 4.69%  | 0.247% |

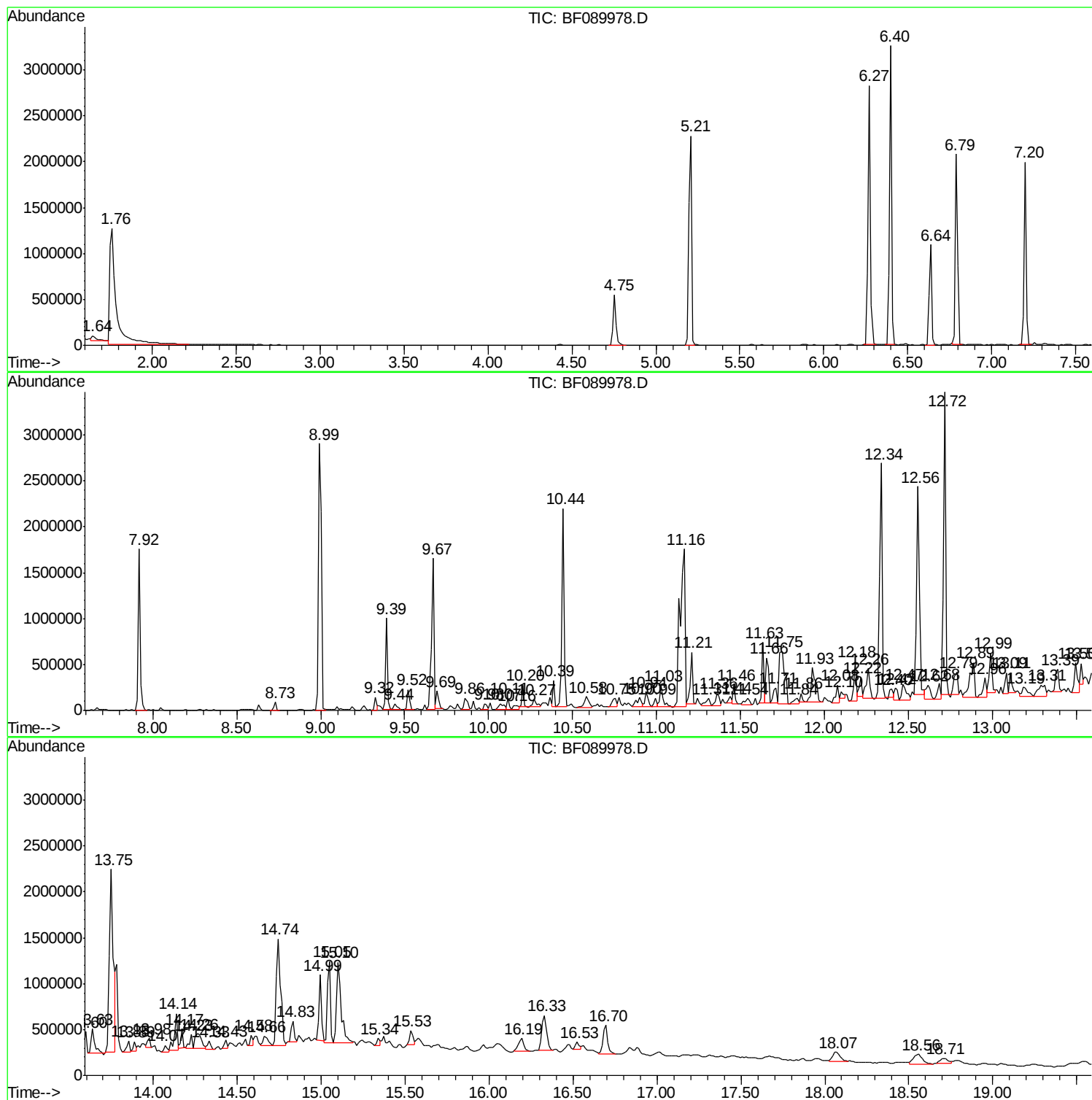
Sum of corrected areas: 69044048

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Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.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\BF090216\  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

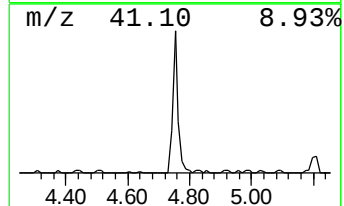
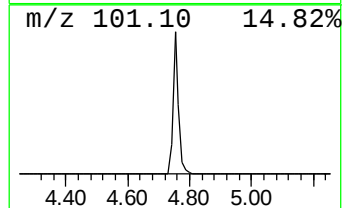
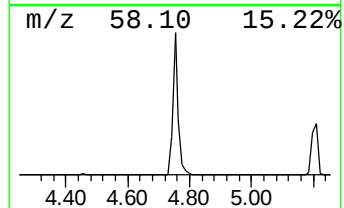
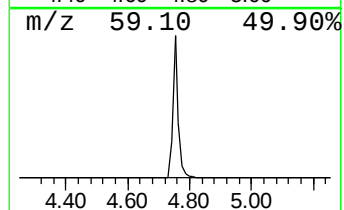
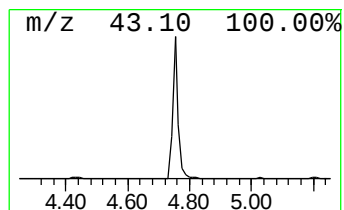
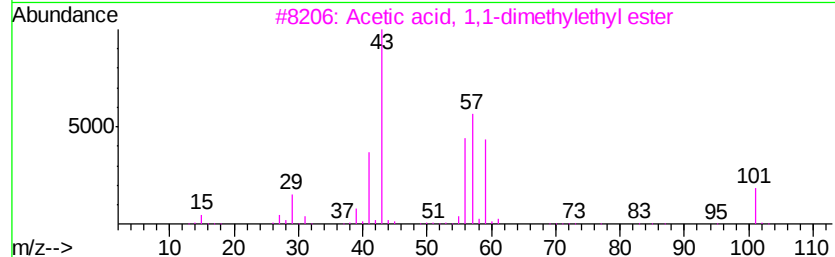
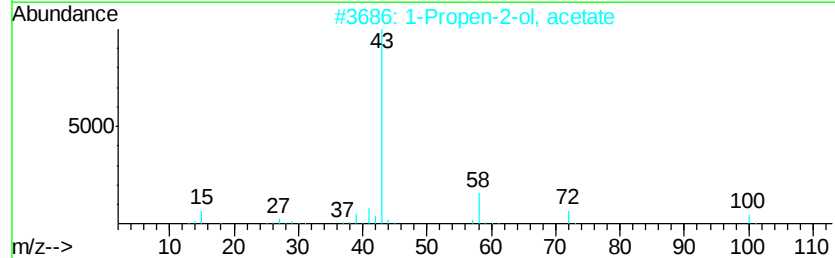
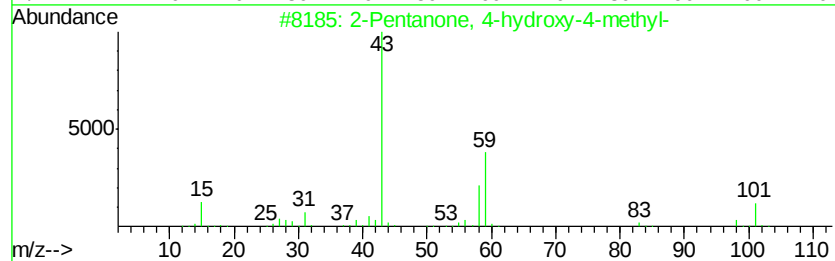
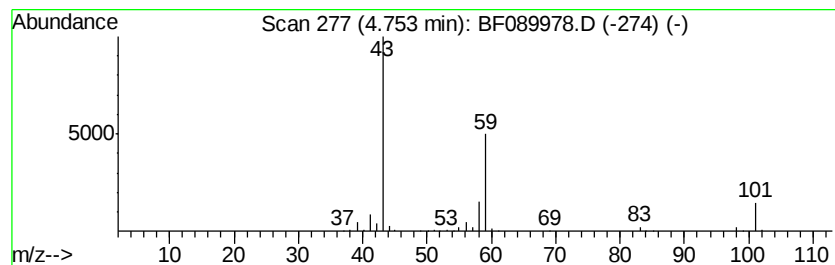
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.75 | 11.54 ng | 673286 | 1,4-Dichlorobenzene-d4 | 6.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2 | 59   |
| 2    |    | 1-Propen-2-ol, acetate              | 100 | C5H8O2  | 000108-22-5 | 10   |
| 3    |    | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 9    |
| 4    |    | Formamide, N,N-diethyl-             | 101 | C5H11NO | 000617-84-5 | 7    |
| 5    |    | (+)-4-Amino-4,5-dihydro-2(3H)-f...  | 101 | C4H7NO2 | 016504-58-8 | 7    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

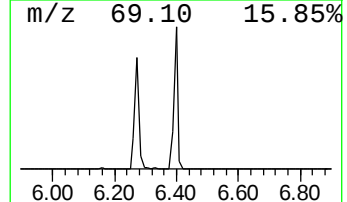
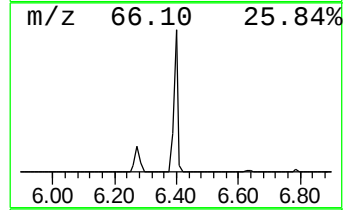
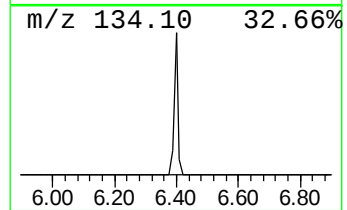
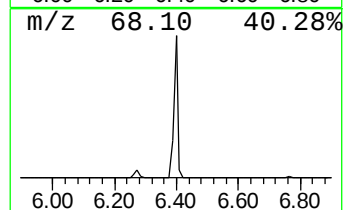
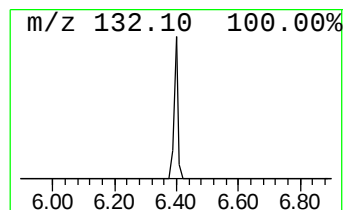
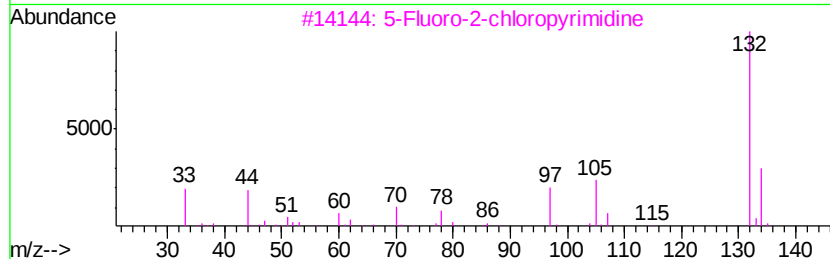
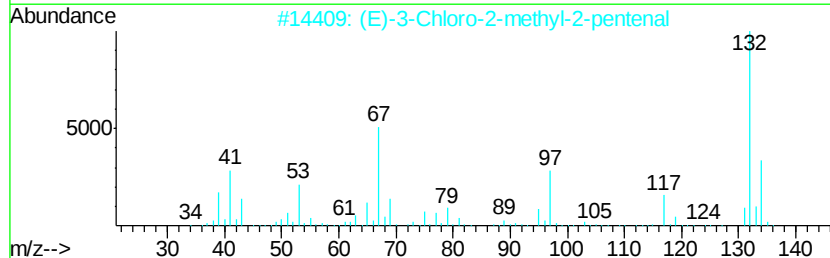
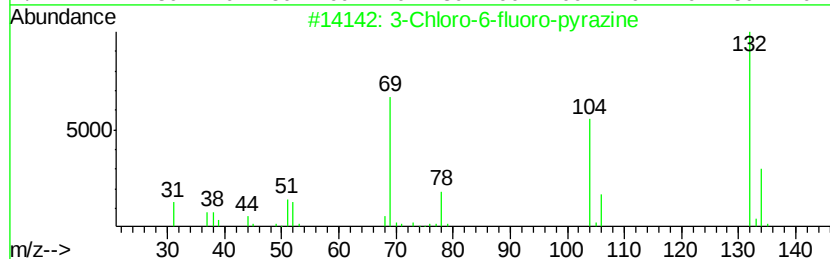
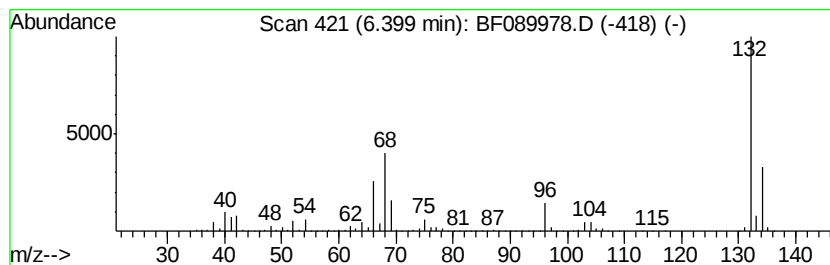
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown6.40 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.40 | 50.34 ng | 2936370 | 1,4-Dichlorobenzene-d4 | 6.64 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 16   |
| 4    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 5    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
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Operator : UM/SJ  
Sample : H4683-01  
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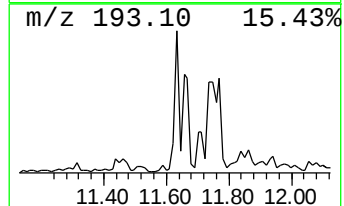
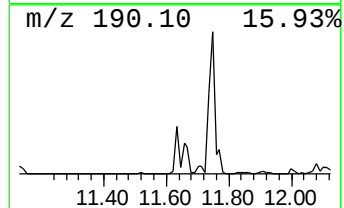
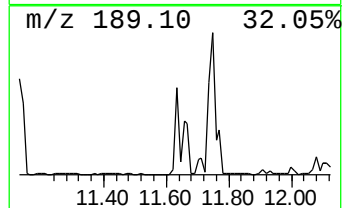
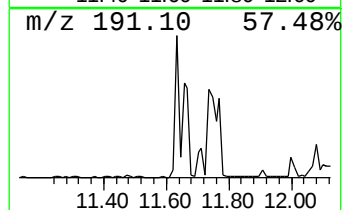
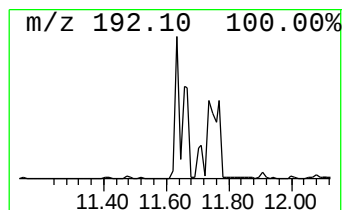
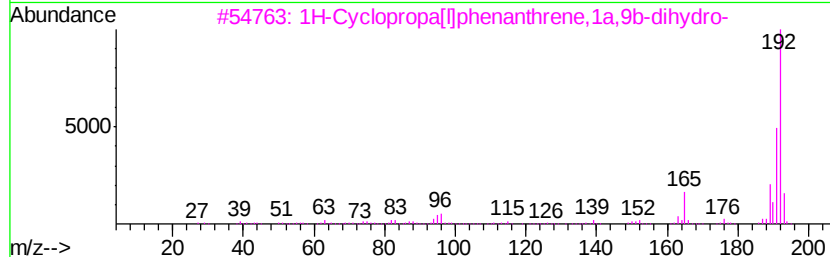
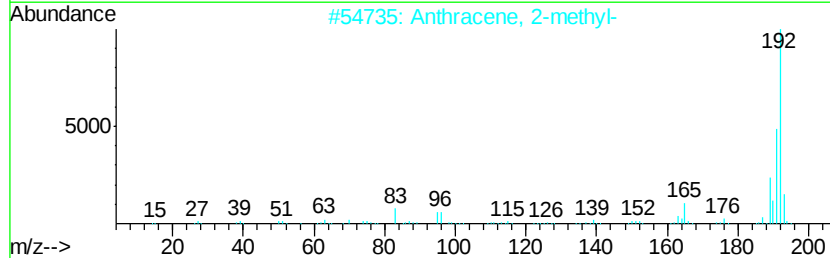
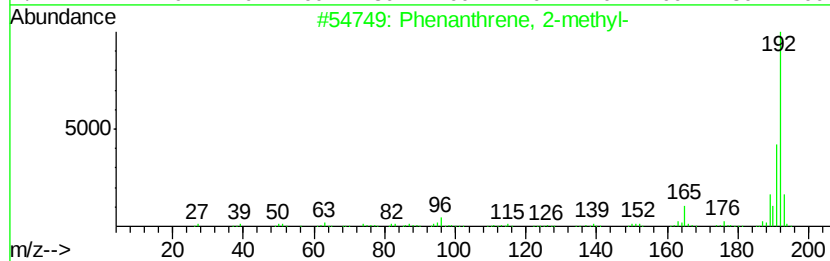
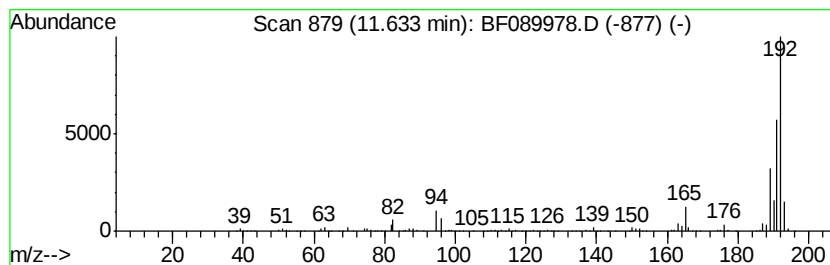
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 14

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.63     | 3.07 ng                             | 559012 | Phenanthrene-d10 | 11.13       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 96   |
| 2         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 96   |
| 3         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 95   |
| 4         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 93   |
| 5         | Anthracene, 9-methyl-               | 192    | C15H12           | 000779-02-2 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
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Operator : UM/SJ  
Sample : H4683-01  
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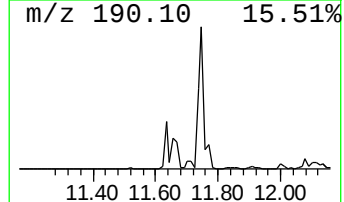
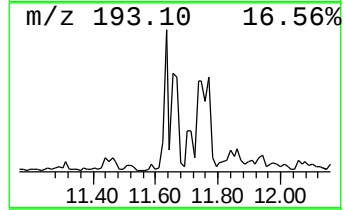
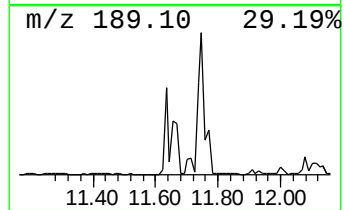
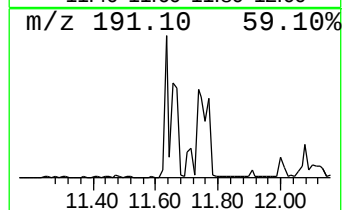
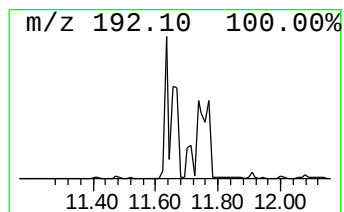
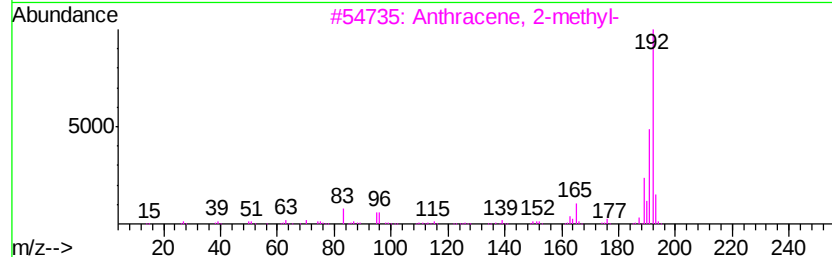
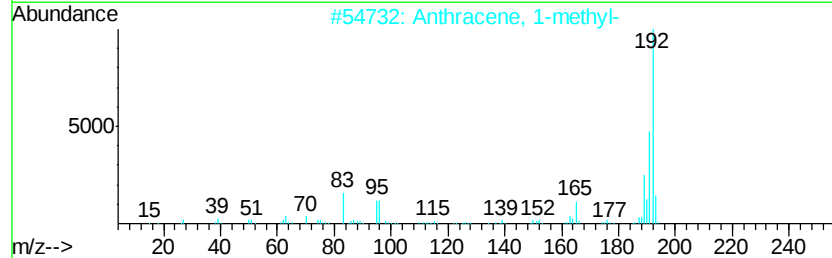
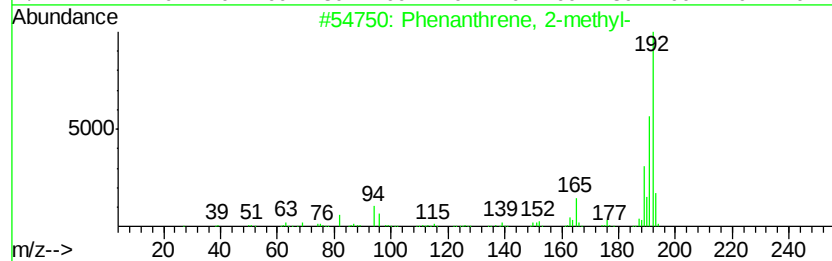
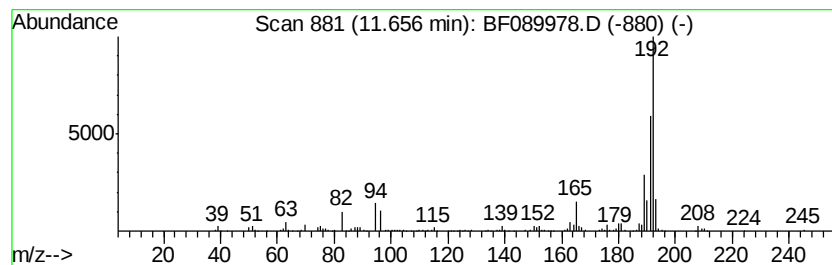
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 Anthracene, 1-methyl- Concentration Rank 13

| R.T.      | EstConc                               | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------------------------|--------|------------------|-------------|------|
| 11.66     | 3.27 ng                               | 594699 | Phenanthrene-d10 | 11.13       |      |
| Hit# of 5 | Tentative ID                          | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2-methyl-               | 192    | C15H12           | 002531-84-2 | 98   |
| 2         | Anthracene, 1-methyl-                 | 192    | C15H12           | 000610-48-0 | 96   |
| 3         | Anthracene, 2-methyl-                 | 192    | C15H12           | 000613-12-7 | 96   |
| 4         | Phenanthrene, 1-methyl-               | 192    | C15H12           | 000832-69-9 | 96   |
| 5         | 1H-Cyclopropa[1]phenanthrene, 1a, ... | 192    | C15H12           | 000949-41-7 | 95   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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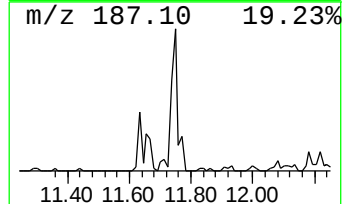
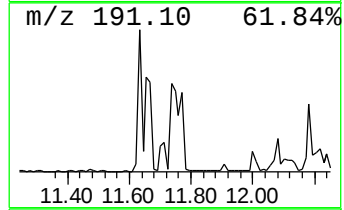
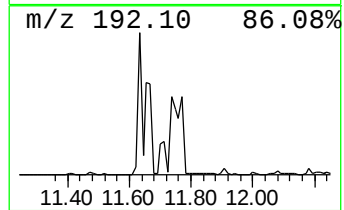
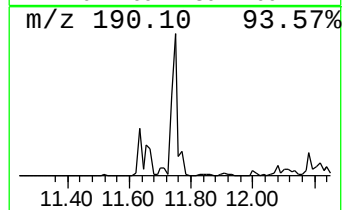
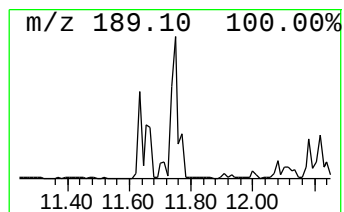
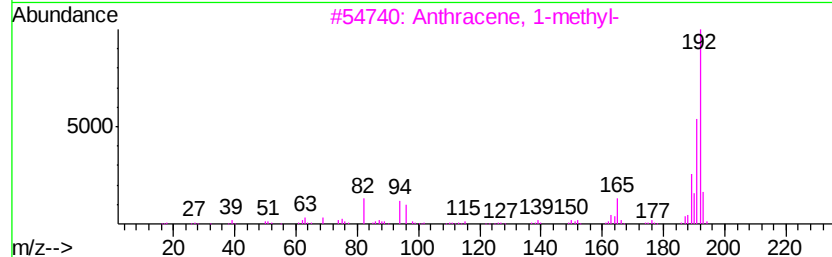
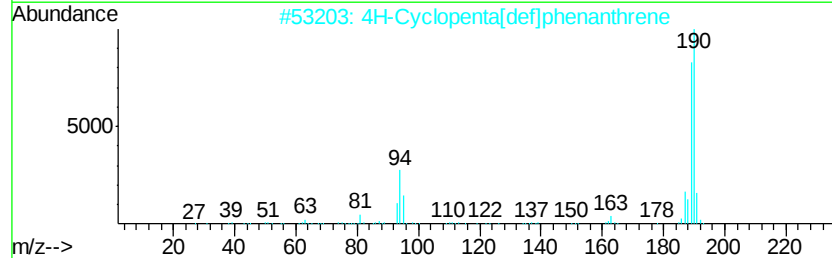
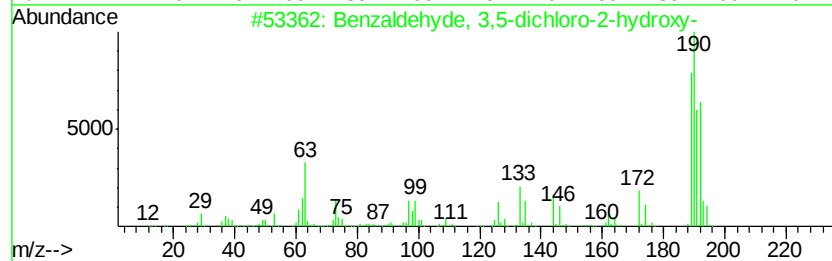
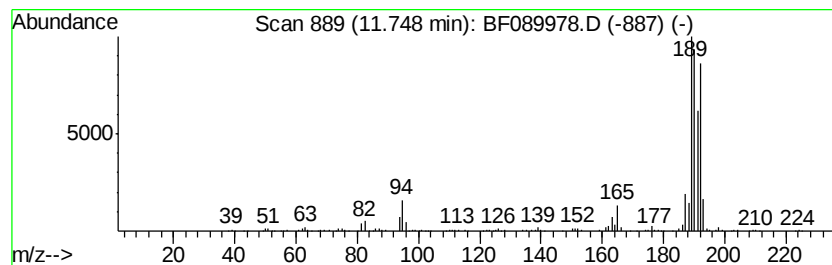
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Benzaldehyde, 3,5-dichloro-... Concentration Rank 6

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 11.75 | 6.68 ng | 1216470 | Phenanthrene-d10 | 11.13 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|---------|---|-------------------------------------|-----|-----------|-------------|------|
| 1       |   | Benzaldehyde, 3,5-dichloro-2-hyd... | 190 | C7H4Cl2O2 | 000090-60-8 | 72   |
| 2       |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10    | 000203-64-5 | 50   |
| 3       |   | Anthracene, 1-methyl-               | 192 | C15H12    | 000610-48-0 | 46   |
| 4       |   | 1H-Indene, 2-phenyl-                | 192 | C15H12    | 004505-48-0 | 42   |
| 5       |   | Anthracene, 9-methyl-               | 192 | C15H12    | 000779-02-2 | 41   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
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Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

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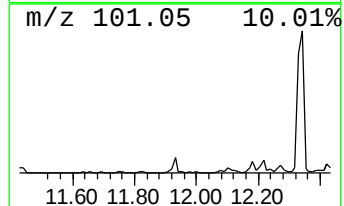
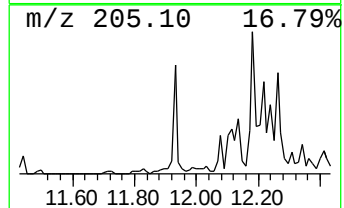
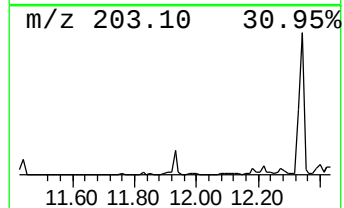
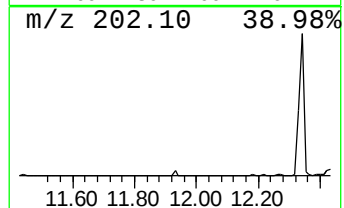
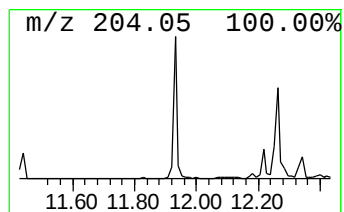
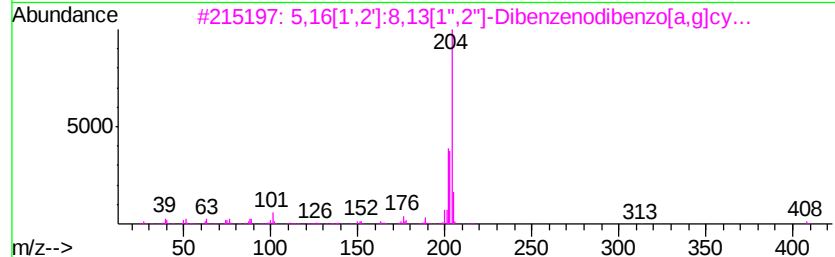
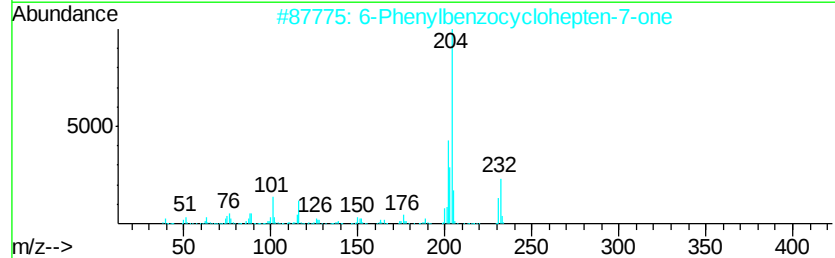
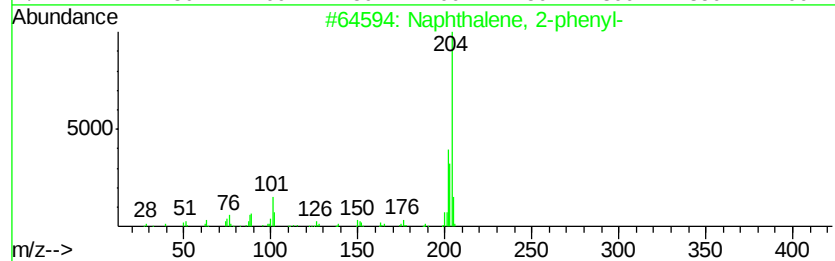
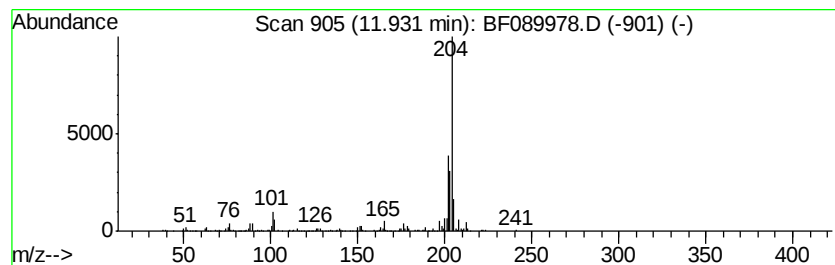
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Naphthalene, 2-phenyl- Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.93 | 3.65 ng | 664984 | Phenanthrene-d10 | 11.13 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 95   |
| 2    |    | 6-Phenylbenzocyclohepten-7-one      | 232 | C17H12O | 093327-56-1 | 90   |
| 3    |    | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24  | 005672-97-9 | 72   |
| 4    |    | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12  | 006572-60-7 | 70   |
| 5    |    | Naphthalene, 1-phenyl-              | 204 | C16H12  | 000605-02-7 | 55   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
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Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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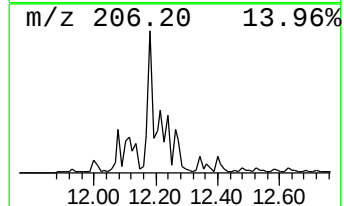
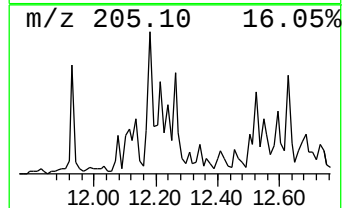
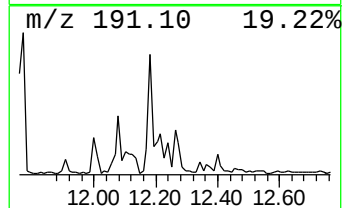
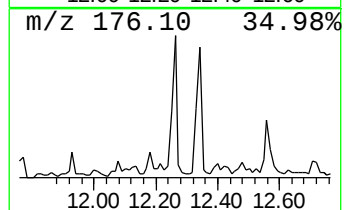
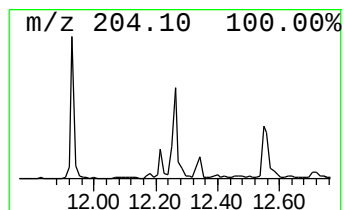
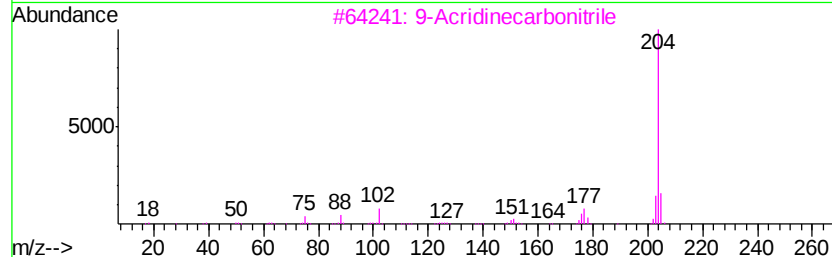
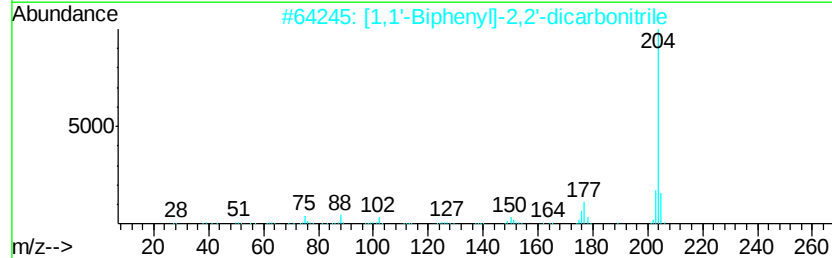
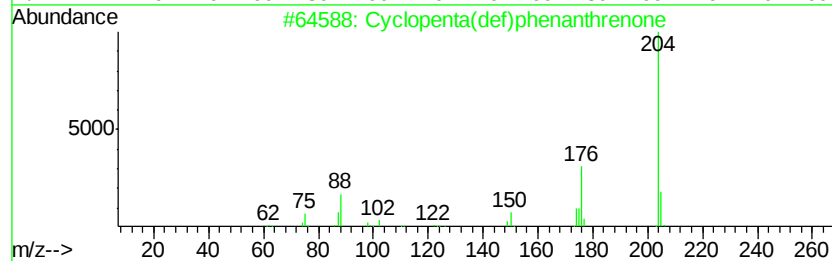
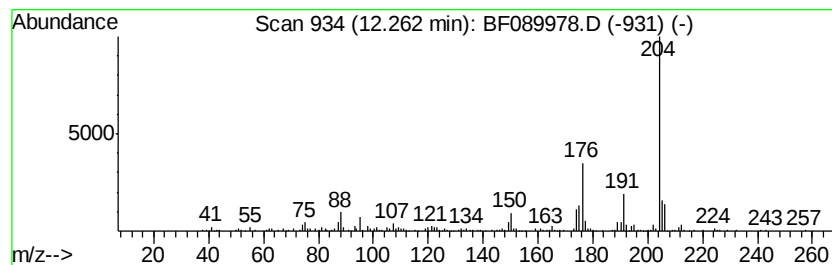
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Cyclopenta(def)phenanthrenone Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.26 | 2.65 ng | 482755 | Phenanthrene-d10 | 11.13 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrenone       | 204 | C15H8O    | 005737-13-3 | 93   |
| 2    |    | [1,1'-Biphenyl]-2,2'-dicarbonitrile | 204 | C14H8N2   | 004341-02-0 | 47   |
| 3    |    | 9-Acridinecarbonitrile              | 204 | C14H8N2   | 005326-19-2 | 47   |
| 4    |    | Quinoline, 6-methoxy-8-nitro-       | 204 | C10H8N2O3 | 000085-81-4 | 47   |
| 5    |    | 1H-3a,7-Methanoazulen-6-ol, octa... | 264 | C17H28O2  | 000077-54-3 | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
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Operator : UM/SJ  
Sample : H4683-01  
Misc :  
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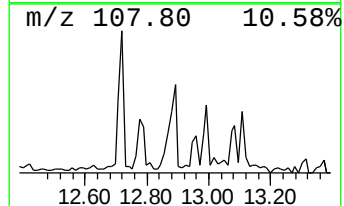
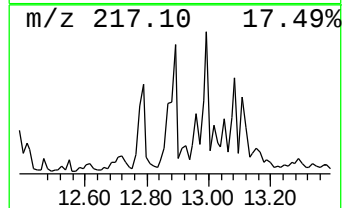
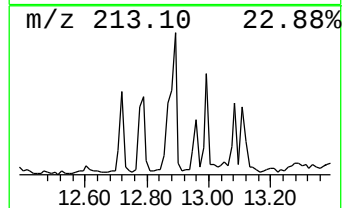
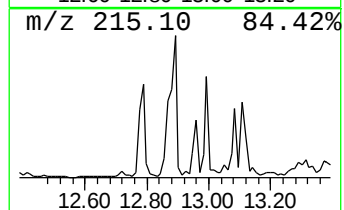
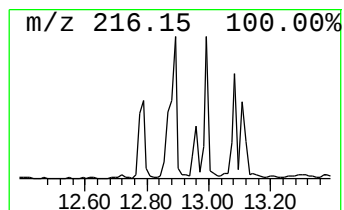
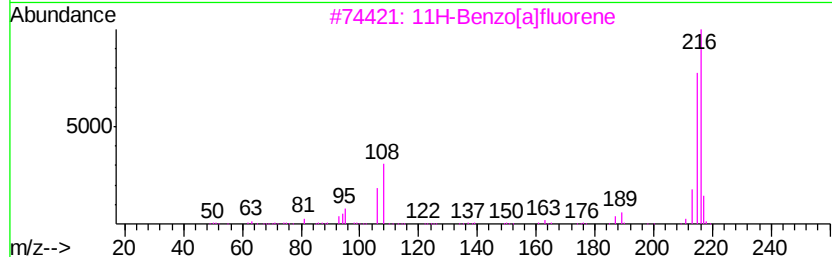
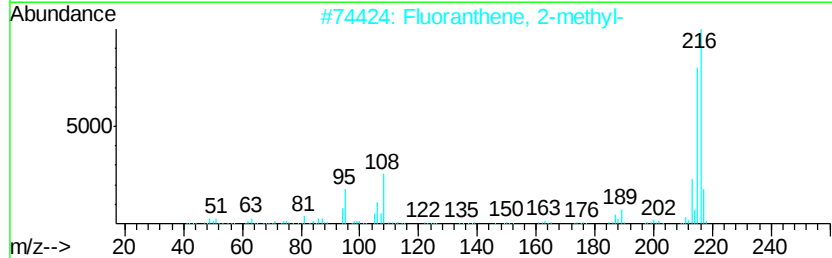
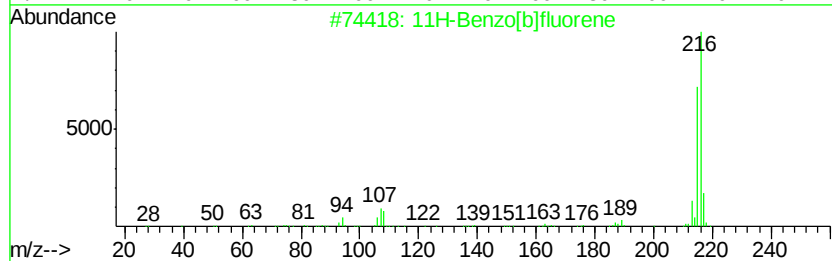
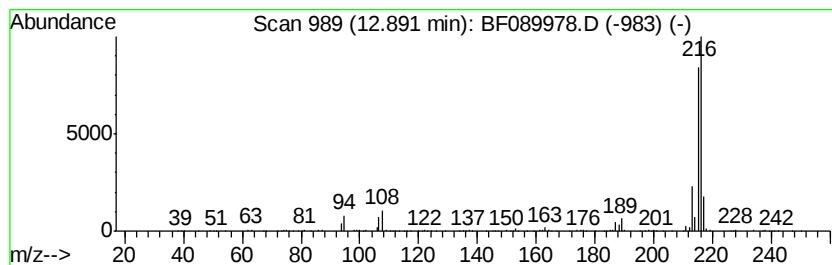
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ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 8

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.    |             |      |
|---------|---------|-------------------------|------------------|---------|-------------|------|
| 12.89   | 4.47 ng | 766495                  | Chrysene-d12     | 13.75   |             |      |
| Hit# of | 5       | Tentative ID            | MW               | MolForm | CAS#        | Qual |
| 1       |         | 11H-Benzo[b]fluorene    | 216              | C17H12  | 000243-17-4 | 95   |
| 2       |         | Fluoranthene, 2-methyl- | 216              | C17H12  | 033543-31-6 | 91   |
| 3       |         | 11H-Benzo[a]fluorene    | 216              | C17H12  | 000238-84-6 | 91   |
| 4       |         | Pyrene, 1-methyl-       | 216              | C17H12  | 002381-21-7 | 90   |
| 5       |         | 7H-Benzo[c]fluorene     | 216              | C17H12  | 000205-12-9 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :  
SOILSAMPLE-1

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

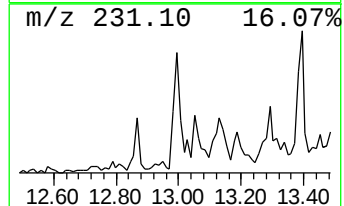
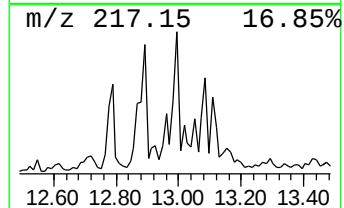
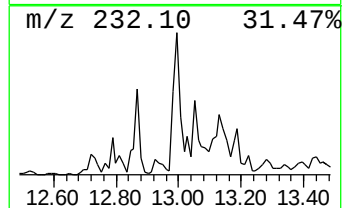
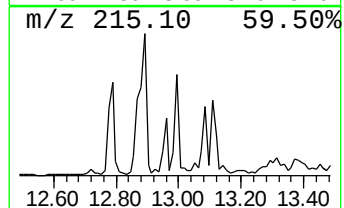
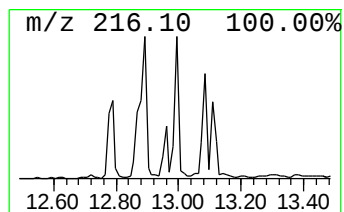
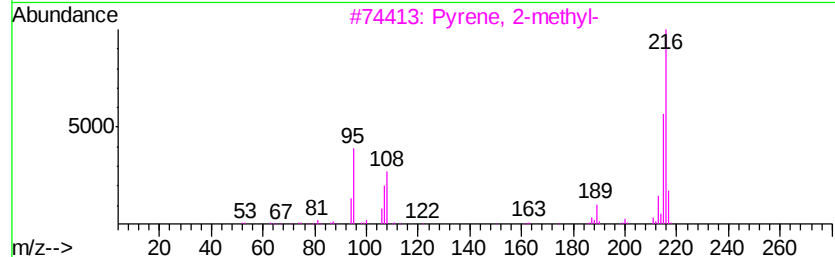
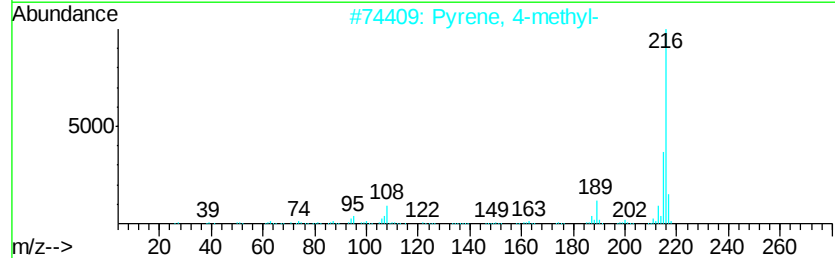
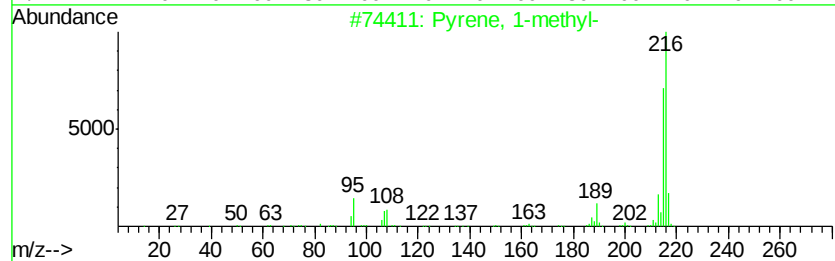
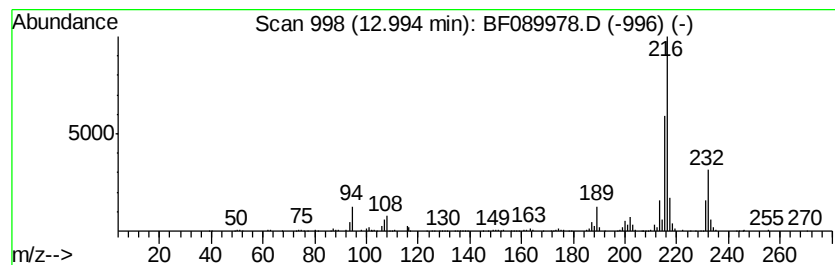
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Pyrene, 1-methyl- Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.99 | 2.58 ng | 442722 | Chrysene-d12     | 13.75 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 96   |
| 2    |    | Pyrene, 4-methyl-       | 216 | C17H12  | 003353-12-6 | 96   |
| 3    |    | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 93   |
| 4    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 86   |
| 5    |    | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 76   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

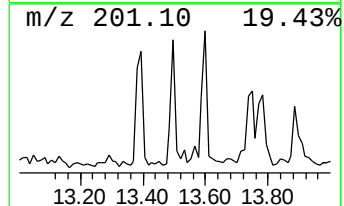
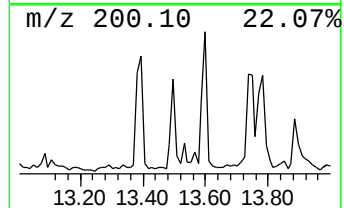
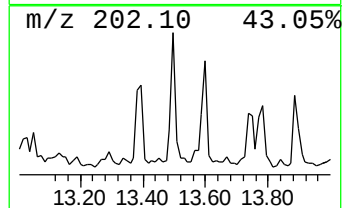
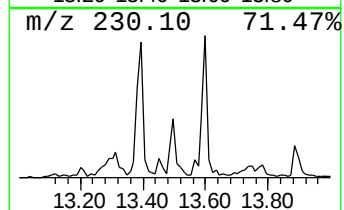
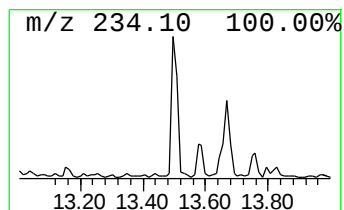
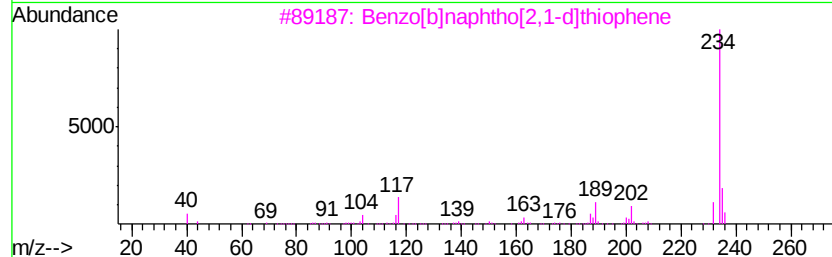
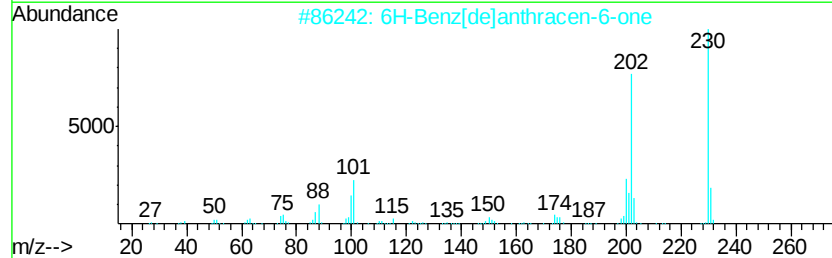
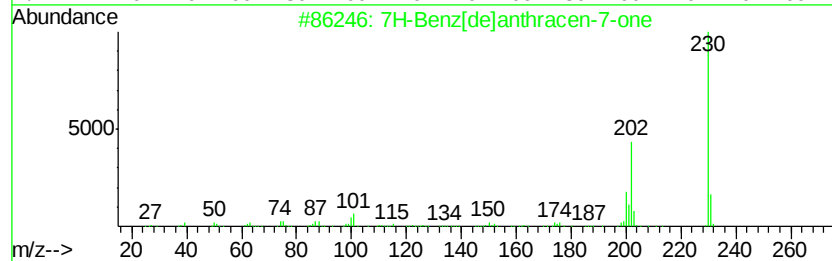
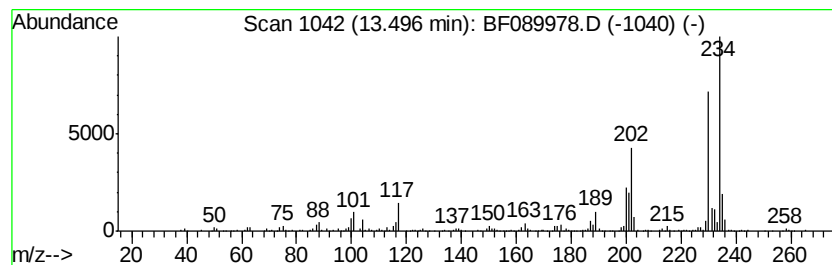
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 7H-Benz[de]anthracen-7-one Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.50 | 2.61 ng | 446981 | Chrysene-d12     | 13.75 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | 7H-Benz[de]anthracen-7-one      | 230 | C17H10O | 000082-05-3 | 90   |
| 2    |    | 6H-Benz[de]anthracen-6-one      | 230 | C17H10O | 080252-14-8 | 86   |
| 3    |    | Benzo[b]naphtho[2,1-d]thiophene | 234 | C16H10S | 000239-35-0 | 86   |
| 4    |    | Anthra(1,2-b)thiophene          | 234 | C16H10S | 000227-86-1 | 83   |
| 5    |    | Benzo[b]naphtho[2,3-d]thiophene | 234 | C16H10S | 000243-46-9 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

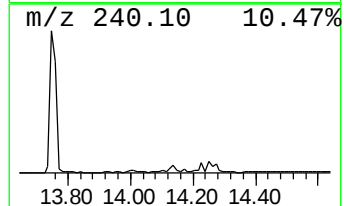
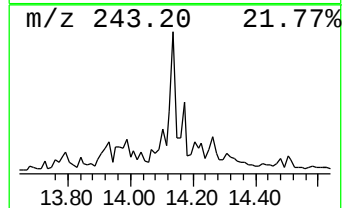
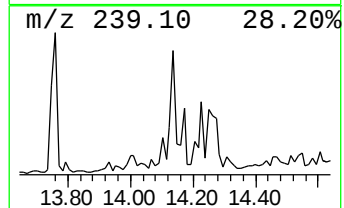
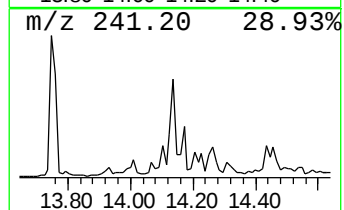
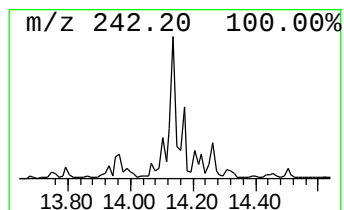
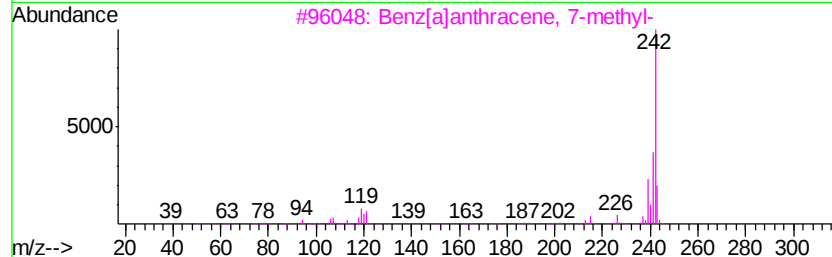
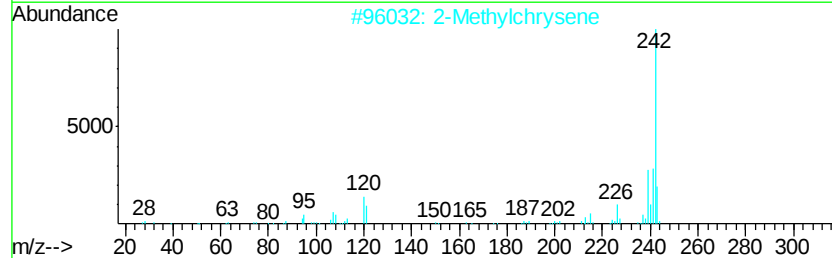
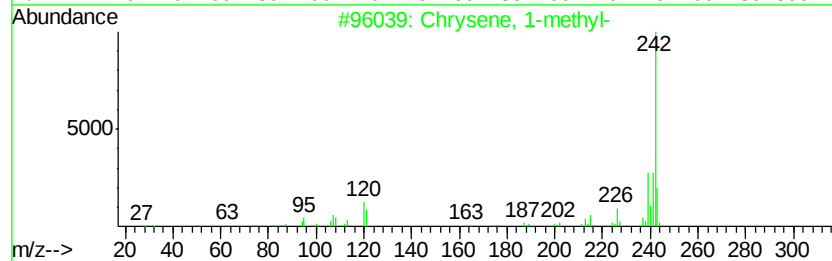
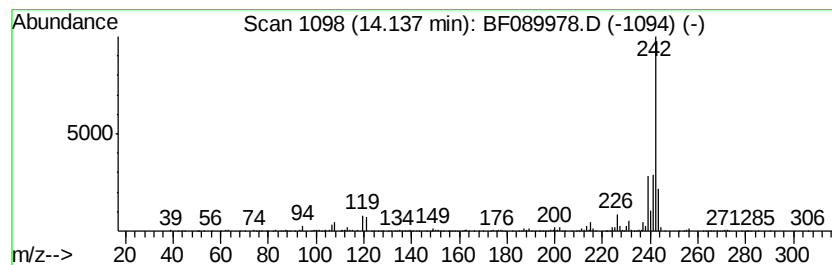
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Chrysene, 1-methyl- Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.14 | 2.87 ng | 492759 | Chrysene-d12     | 13.75 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Chrysene, 1-methyl-          | 242 | C19H14  | 003351-28-8 | 98   |
| 2    |    | 2-Methylchrysene             | 242 | C19H14  | 003351-32-4 | 96   |
| 3    |    | Benz[a]anthracene, 7-methyl- | 242 | C19H14  | 002541-69-7 | 96   |
| 4    |    | Triphenylene, 2-methyl-      | 242 | C19H14  | 001705-84-6 | 96   |
| 5    |    | Benz[a]anthracene, 8-methyl- | 242 | C19H14  | 002381-31-9 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

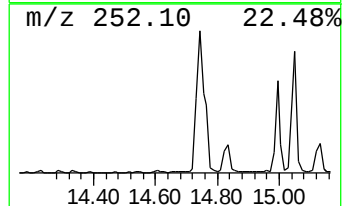
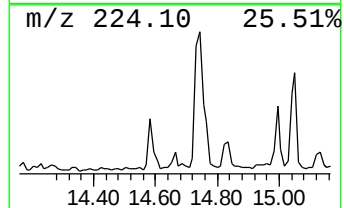
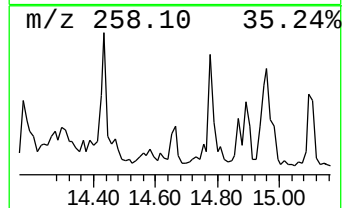
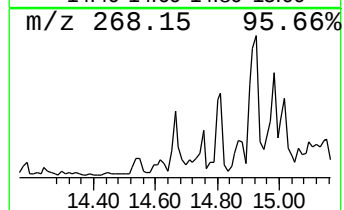
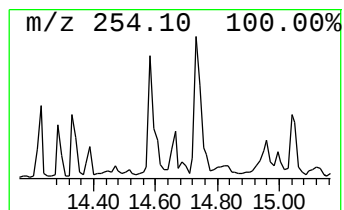
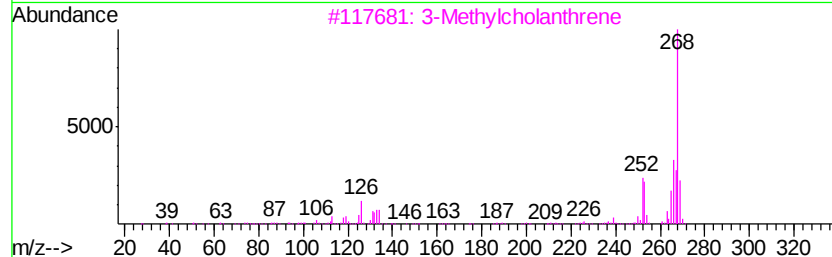
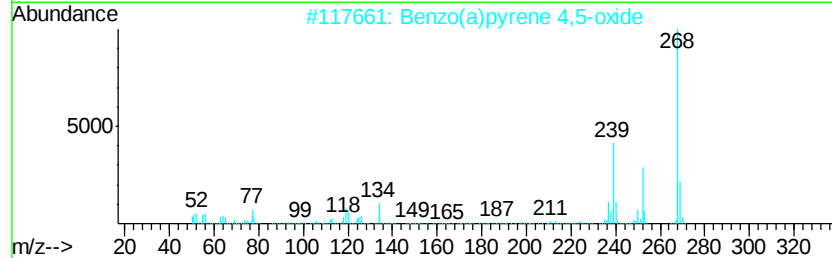
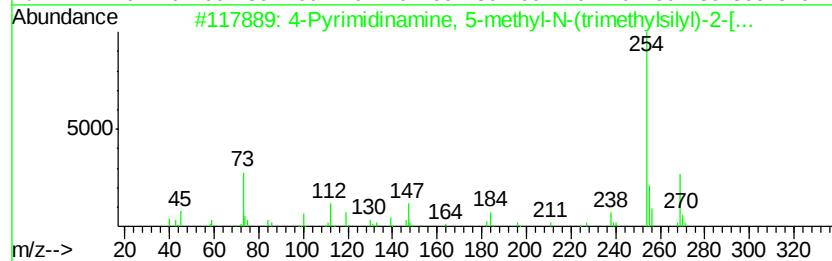
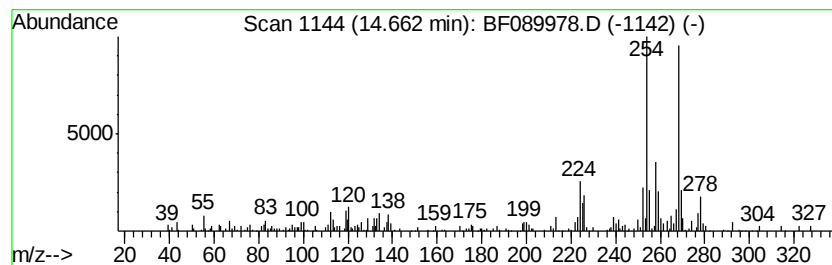
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 unknown14.66 Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.66 | 2.58 ng | 199969 | Perylene-d12     | 15.10 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|--------------|--------------|------|
| 1    |    |   | 4-Pyrimidinamine, 5-methyl-N-(tr... | 269 | C11H23N3OSi2 | 032865-88-6  | 42   |
| 2    |    |   | Benzo(a)pyrene 4,5-oxide            | 268 | C20H12O      | 037574-47-3  | 38   |
| 3    |    |   | 3-Methylcholanthrene                | 268 | C21H16       | 000056-49-5  | 30   |
| 4    |    |   | trans-4'-Methyl-4-(methylthio)ch... | 268 | C17H16OS     | 126443-27-4  | 30   |
| 5    |    |   | 2-Cyclohexen-1-one, 3-[3-(4,4-di... | 269 | C17H19NO2    | 1000197-92-1 | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

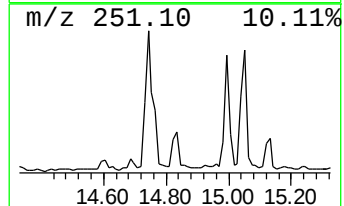
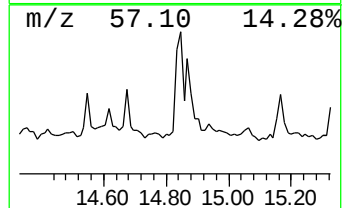
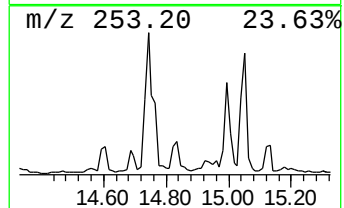
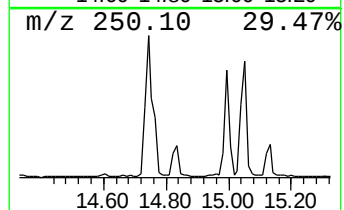
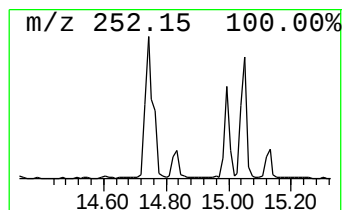
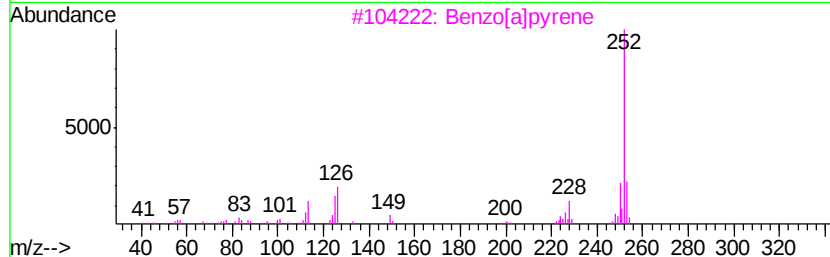
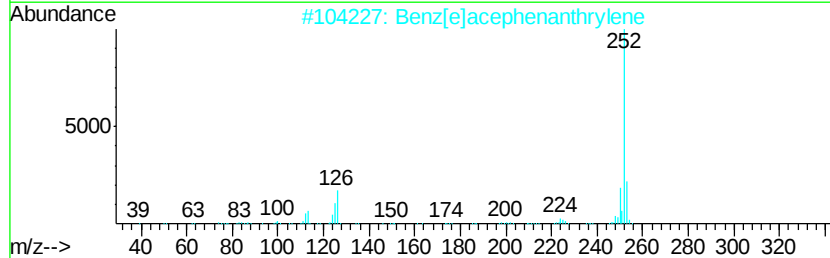
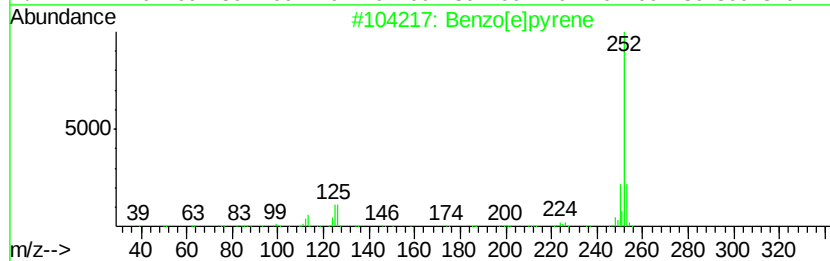
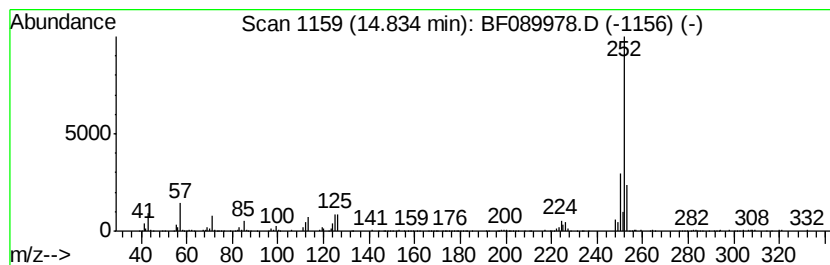
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzo[e]pyrene Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.83 | 3.78 ng | 292420 | Perylene-d12     | 15.10 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[e]pyrene           | 252 | C20H12  | 000192-97-2 | 94   |
| 2    |    |   | Benz[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 87   |
| 3    |    |   | Benzo[a]pyrene           | 252 | C20H12  | 000050-32-8 | 87   |
| 4    |    |   | Perylene                 | 252 | C20H12  | 000198-55-0 | 76   |
| 5    |    |   | Benzo[k]fluoranthene     | 252 | C20H12  | 000207-08-9 | 76   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

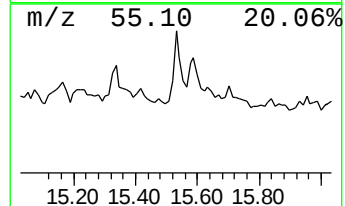
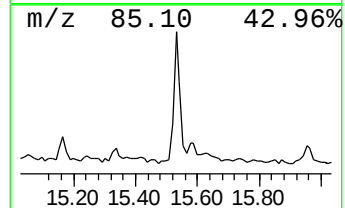
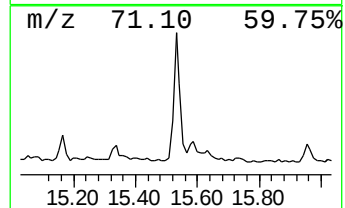
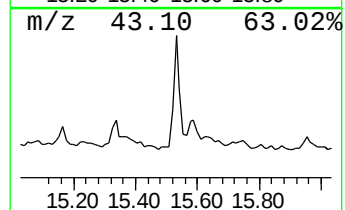
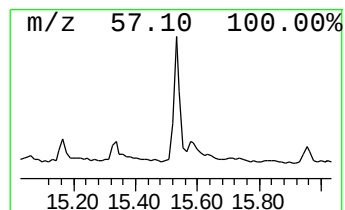
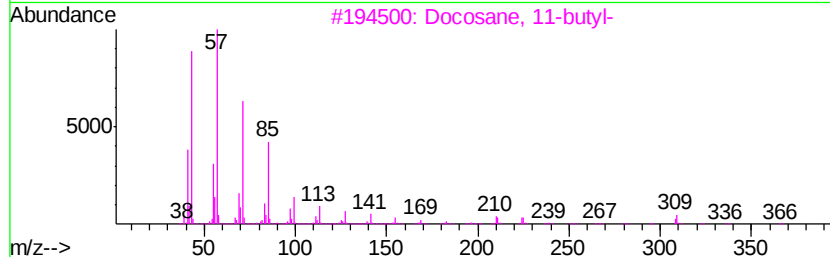
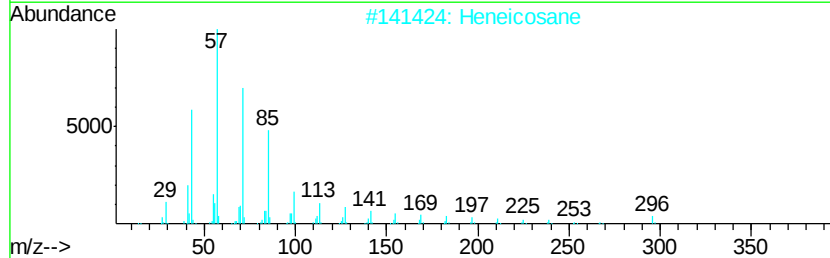
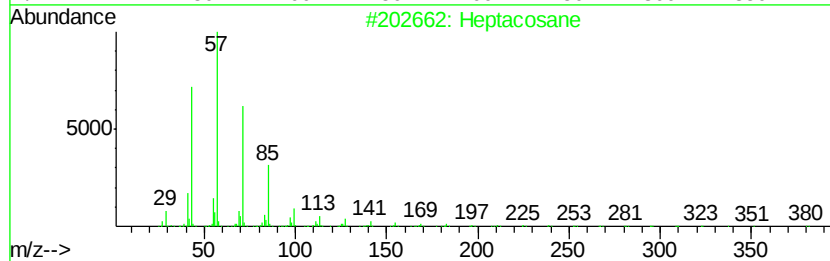
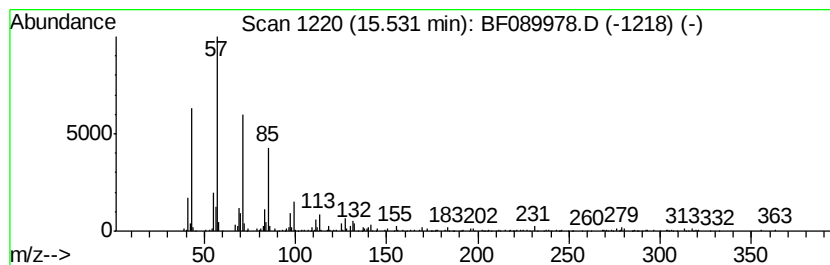
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Heptacosane Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.53 | 3.00 ng | 232164 | Perylene-d12     | 15.10 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Heptacosane         | 380 | C27H56  | 000593-49-7 | 90   |
| 2    |    | Heneicosane         | 296 | C21H44  | 000629-94-7 | 87   |
| 3    |    | Docosane, 11-butyl- | 366 | C26H54  | 013475-76-8 | 86   |
| 4    |    | Eicosane, 7-hexyl-  | 366 | C26H54  | 055333-99-8 | 86   |
| 5    |    | Heptadecane         | 240 | C17H36  | 000629-78-7 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

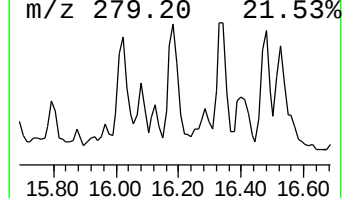
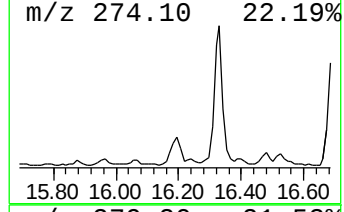
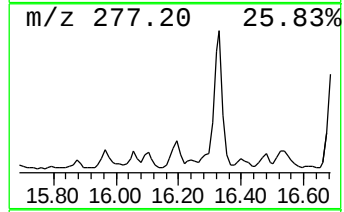
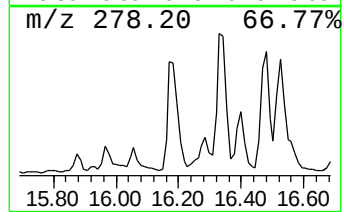
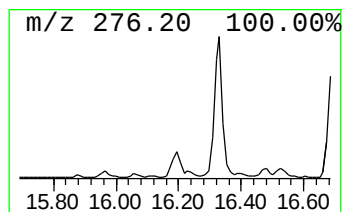
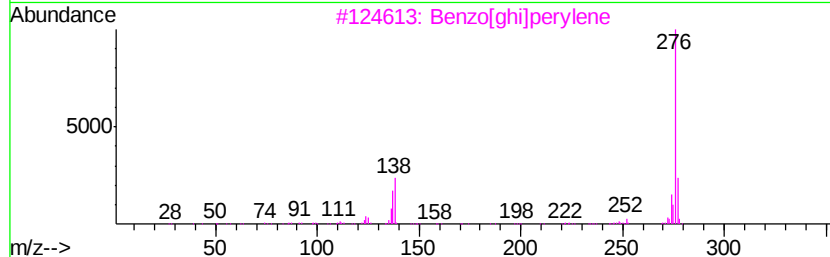
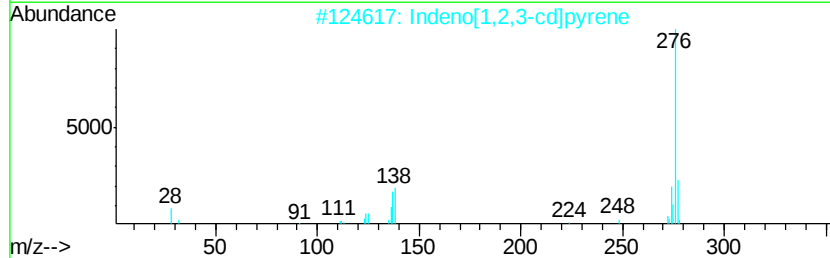
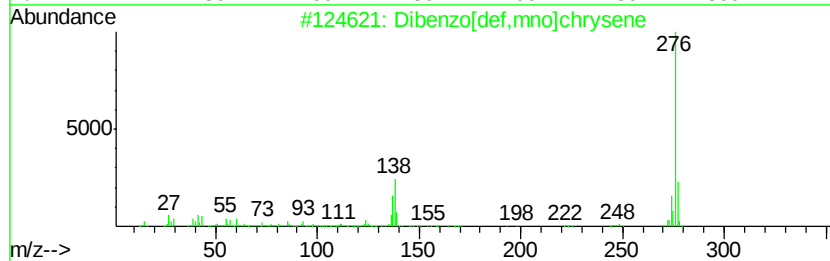
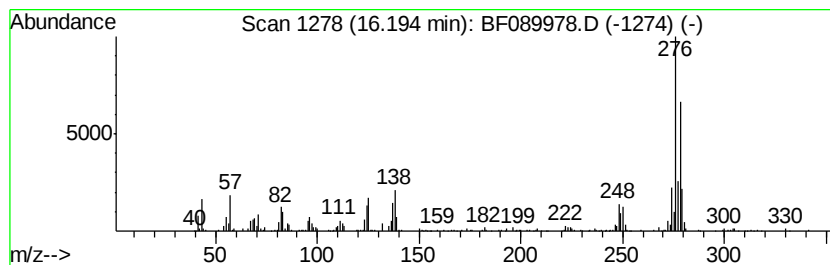
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Dibenzo[def,mno]chrysene Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.19 | 4.21 ng | 326272 | Perylene-d12     | 15.10 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Dibenzo[def,mno]chrysene            | 276 | C22H12   | 000191-26-4 | 87   |
| 2    |    | Indeno[1,2,3-cd]pyrene              | 276 | C22H12   | 000193-39-5 | 70   |
| 3    |    | Benzo[ghi]perylene                  | 276 | C22H12   | 000191-24-2 | 55   |
| 4    |    | Indeno[1,2,3-cd]fluoranthene        | 276 | C22H12   | 000193-43-1 | 38   |
| 5    |    | 1-Naphthalenecarboxylic acid, 2-... | 276 | C18H12O3 | 038119-11-8 | 30   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

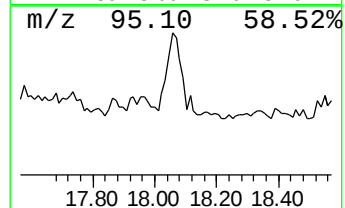
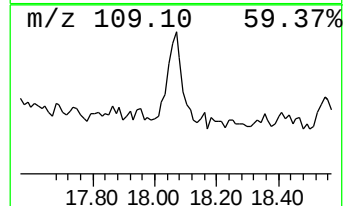
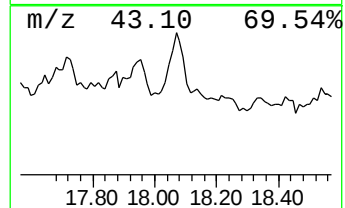
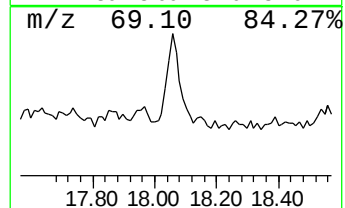
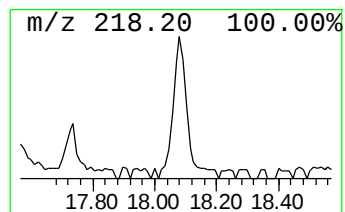
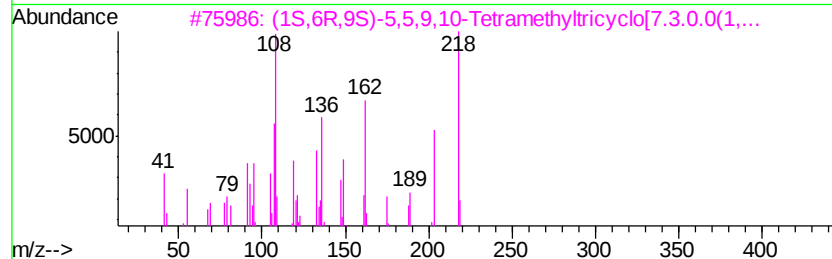
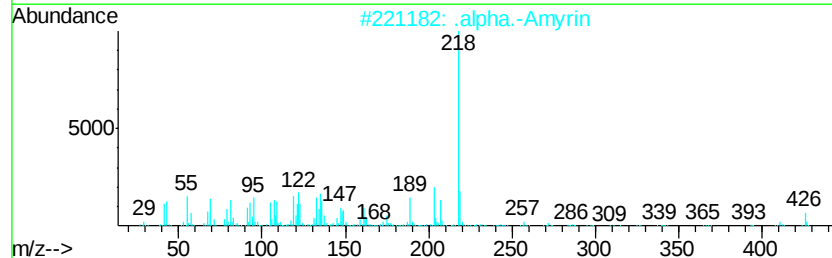
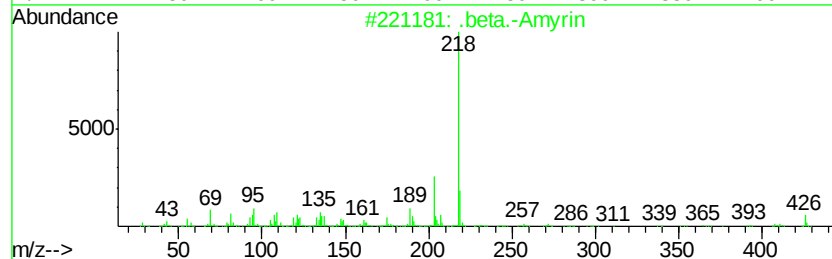
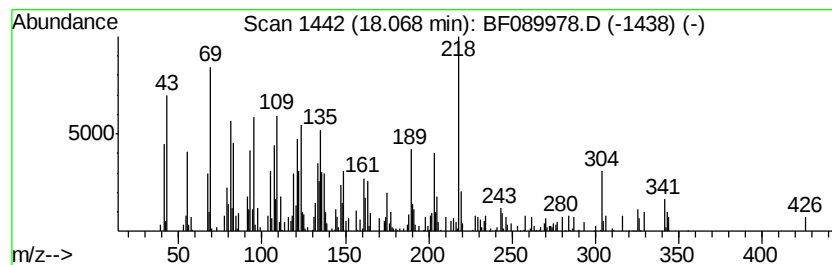
Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 .beta.-Amyrin Concentration Rank 10

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.            |
|---------|---------|-------------------------------------|------------------|-----------------|
| 18.07   | 3.79 ng | 293079                              | Perylene-d12     | 15.10           |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |         | .beta.-Amyrin                       | 426 C30H500      | 000559-70-6 80  |
| 2       |         | .alpha.-Amyrin                      | 426 C30H500      | 000638-95-9 58  |
| 3       |         | (1S,6R,9S)-5,5,9,10-Tetramethylt... | 218 C16H26       | 1000298-97-8 52 |
| 4       |         | 9-Isopropyl-6-methyl-6,7-dihydro... | 219 C13H17N02    | 134076-78-1 38  |
| 5       |         | 5(1H)-Azulenone, 2,4,6,7,8,8a-he... | 218 C15H220      | 006754-66-1 35  |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
Data File : BF089978.D  
Acq On : 2 Sep 2016 16:43  
Operator : UM/SJ  
Sample : H4683-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

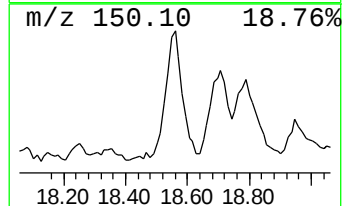
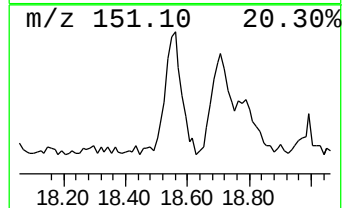
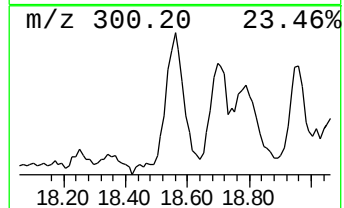
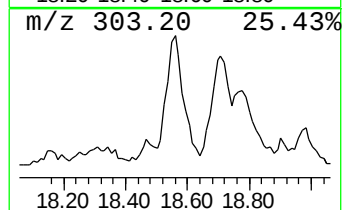
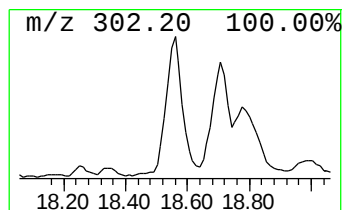
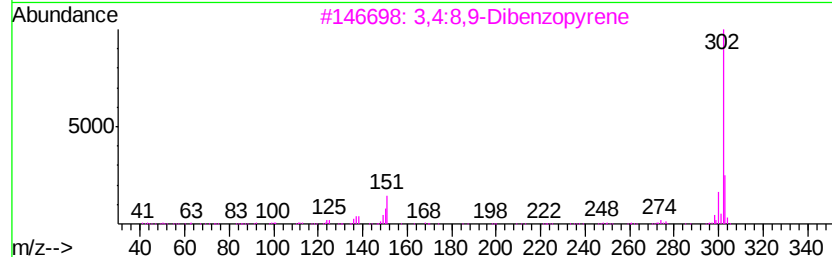
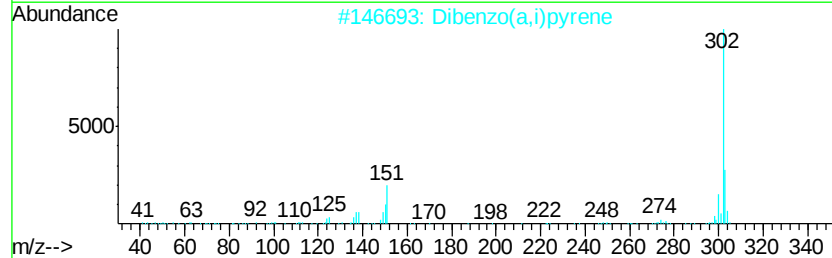
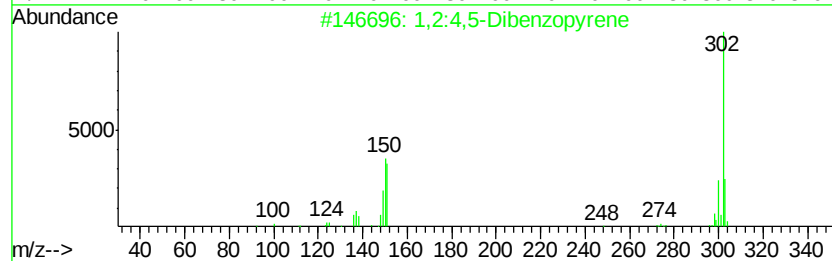
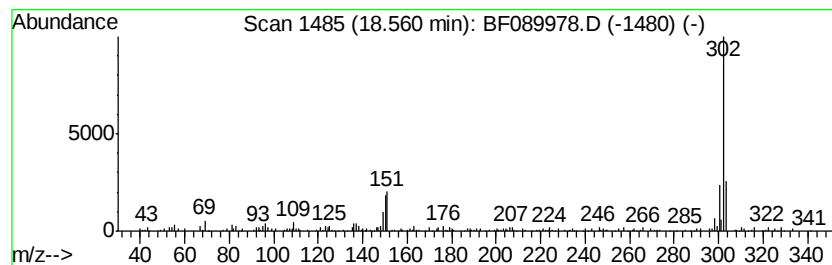
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 1,2:4,5-Dibenzopyrene Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.56 | 5.31 ng | 410924 | Perylene-d12     | 15.10 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | 1,2:4,5-Dibenzopyrene     | 302 | C24H14  | 000192-65-4 | 93   |
| 2    |    | Dibenzo(a,i)pyrene        | 302 | C24H14  | 000189-55-9 | 90   |
| 3    |    | 3,4:8,9-Dibenzopyrene     | 302 | C24H14  | 000189-64-0 | 90   |
| 4    |    | Dibenzo[fg,op]naphthacene | 302 | C24H14  | 000192-51-8 | 64   |
| 5    |    | Dibenz(a,e)aceanthrylene  | 302 | C24H14  | 005385-75-1 | 58   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF090216\  
 Data File : BF089978.D  
 Acq On : 2 Sep 2016 16:43  
 Operator : UM/SJ  
 Sample : H4683-01  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SOILSAMPLE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF083116.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... | 4.75  | 11.5    | ng    | 673286   | 1                     | 6.64  | 1166710 | 20.0 |
| unknown6.40          | 6.40  | 50.3    | ng    | 2936370  | 1                     | 6.64  | 1166710 | 20.0 |
| Phenanthrene, 2-m... | 11.63 | 3.1     | ng    | 559012   | 4                     | 11.13 | 3641670 | 20.0 |
| Anthracene, 1-met... | 11.66 | 3.3     | ng    | 594699   | 4                     | 11.13 | 3641670 | 20.0 |
| Benzaldehyde, 3,5... | 11.75 | 6.7     | ng    | 1216470  | 4                     | 11.13 | 3641670 | 20.0 |
| Naphthalene, 2-ph... | 11.93 | 3.6     | ng    | 664984   | 4                     | 11.13 | 3641670 | 20.0 |
| Cyclopenta(def)ph... | 12.26 | 2.6     | ng    | 482755   | 4                     | 11.13 | 3641670 | 20.0 |
| 11H-Benzo[b]fluorene | 12.89 | 4.5     | ng    | 766495   | 5                     | 13.75 | 3430970 | 20.0 |
| Pyrene, 1-methyl-    | 12.99 | 2.6     | ng    | 442722   | 5                     | 13.75 | 3430970 | 20.0 |
| 7H-Benz[de]anthra... | 13.50 | 2.6     | ng    | 446981   | 5                     | 13.75 | 3430970 | 20.0 |
| Chrysene, 1-methyl-  | 14.14 | 2.9     | ng    | 492759   | 5                     | 13.75 | 3430970 | 20.0 |
| unknown14.66         | 14.66 | 2.6     | ng    | 199969   | 6                     | 15.10 | 1548510 | 20.0 |
| Benzo[e]pyrene       | 14.83 | 3.8     | ng    | 292420   | 6                     | 15.10 | 1548510 | 20.0 |
| Heptacosane          | 15.53 | 3.0     | ng    | 232164   | 6                     | 15.10 | 1548510 | 20.0 |
| Dibenzo[def,mno]c... | 16.19 | 4.2     | ng    | 326272   | 6                     | 15.10 | 1548510 | 20.0 |
| .beta.-Amyrin        | 18.07 | 3.8     | ng    | 293079   | 6                     | 15.10 | 1548510 | 20.0 |
| 1,2:4,5-Dibenzopy... | 18.56 | 5.3     | ng    | 410924   | 6                     | 15.10 | 1548510 | 20.0 |