

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
 Data File : BF079686.D  
 Acq On : 4 Jun 2015 4:28  
 Operator : TP/IZ  
 Sample : G2408-25  
 Misc :  
 ALS Vial : 32 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-14D

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-BF060315.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         | 1.393       | 17            | 18          | 20           | rVB      | 17473645       | 12093624      | 100.00%         | 15.574%       |
| 2         | 1.518       | 27            | 29          | 49           | rVB3     | 373006         | 1710537       | 14.14%          | 2.203%        |
| 3         | 2.273       | 93            | 95          | 109          | rVB      | 84923          | 273428        | 2.26%           | 0.352%        |
| 4         | 2.478       | 109           | 113         | 116          | rBV      | 4309009        | 7393524       | 61.14%          | 9.522%        |
| 5         | 2.684       | 128           | 131         | 142          | rVB      | 88876          | 200746        | 1.66%           | 0.259%        |
| 6         | 2.947       | 151           | 154         | 158          | rVB      | 259031         | 378025        | 3.13%           | 0.487%        |
| 7         | 6.136       | 430           | 433         | 435          | rBV      | 2605983        | 2215438       | 18.32%          | 2.853%        |
| 8         | 7.108       | 516           | 518         | 521          | rBV      | 2181440        | 2201779       | 18.21%          | 2.835%        |
| 9         | 7.290       | 531           | 534         | 536          | rBV      | 2320148        | 2397411       | 19.82%          | 3.087%        |
| 10        | 7.519       | 552           | 554         | 556          | rBV      | 618864         | 524321        | 4.34%           | 0.675%        |
| 11        | 7.679       | 566           | 568         | 571          | rBV      | 2075966        | 1829478       | 15.13%          | 2.356%        |
| 12        | 8.079       | 600           | 603         | 605          | rBV      | 1075172        | 1276672       | 10.56%          | 1.644%        |
| 13        | 8.811       | 665           | 667         | 672          | rVB      | 716996         | 757991        | 6.27%           | 0.976%        |
| 14        | 9.885       | 759           | 761         | 764          | rVB      | 2641059        | 2773364       | 22.93%          | 3.572%        |
| 15        | 10.274      | 793           | 795         | 797          | rVB      | 377066         | 318658        | 2.63%           | 0.410%        |
| 16        | 10.594      | 821           | 823         | 824          | rBV      | 877352         | 780929        | 6.46%           | 1.006%        |
| 17        | 10.628      | 824           | 826         | 827          | rVB      | 227705         | 226663        | 1.87%           | 0.292%        |
| 18        | 11.142      | 869           | 871         | 873          | rVB      | 244453         | 230186        | 1.90%           | 0.296%        |
| 19        | 11.211      | 873           | 877         | 880          | rBV3     | 99840          | 203895        | 1.69%           | 0.263%        |
| 20        | 11.382      | 890           | 892         | 894          | rBV2     | 1384593        | 1488108       | 12.30%          | 1.916%        |
| 21        | 11.702      | 914           | 920         | 921          | rBV2     | 97202          | 198778        | 1.64%           | 0.256%        |
| 22        | 11.999      | 944           | 946         | 951          | rVB      | 172744         | 179557        | 1.48%           | 0.231%        |
| 23        | 12.102      | 953           | 955         | 956          | rBV      | 923673         | 862932        | 7.14%           | 1.111%        |
| 24        | 12.125      | 956           | 957         | 959          | rVB      | 2247538        | 2288807       | 18.93%          | 2.948%        |
| 25        | 12.182      | 959           | 962         | 964          | rBV      | 627463         | 570683        | 4.72%           | 0.735%        |
| 26        | 12.617      | 999           | 1000        | 1002         | rBV      | 364982         | 478297        | 3.95%           | 0.616%        |
| 27        | 12.651      | 1002          | 1003        | 1005         | rVB      | 722014         | 536105        | 4.43%           | 0.690%        |
| 28        | 12.697      | 1005          | 1007        | 1009         | rBV      | 214248         | 241448        | 2.00%           | 0.311%        |
| 29        | 12.742      | 1009          | 1011        | 1015         | rBV2     | 618368         | 1209209       | 10.00%          | 1.557%        |
| 30        | 12.925      | 1025          | 1027        | 1028         | rBV      | 330144         | 306363        | 2.53%           | 0.395%        |
| 31        | 12.948      | 1028          | 1029        | 1033         | rVB2     | 167319         | 256843        | 2.12%           | 0.331%        |
| 32        | 13.085      | 1036          | 1041        | 1043         | rBV2     | 144740         | 237990        | 1.97%           | 0.306%        |
| 33        | 13.211      | 1049          | 1052        | 1053         | rBV      | 289510         | 430468        | 3.56%           | 0.554%        |
| 34        | 13.245      | 1053          | 1055        | 1058         | rVB2     | 160907         | 356094        | 2.94%           | 0.459%        |

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
 Data File : BF079686.D  
 Acq On : 4 Jun 2015 4:28  
 Operator : TP/IZ  
 Sample : G2408-25  
 Misc :  
 ALS Vial : 32 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-14D

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-BF060315.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 13.302 | 1058 | 1060 | 1065 | rBV2 | 242675  | 365644  | 3.02%  | 0.471% |
| 36 | 13.394 | 1065 | 1068 | 1070 | rBV  | 3046951 | 3151659 | 26.06% | 4.059% |
| 37 | 13.554 | 1080 | 1082 | 1086 | rVB3 | 130693  | 256938  | 2.12%  | 0.331% |
| 38 | 13.645 | 1086 | 1090 | 1093 | rBV  | 2839375 | 3720649 | 30.77% | 4.792% |
| 39 | 13.691 | 1093 | 1094 | 1097 | rVB2 | 190413  | 224124  | 1.85%  | 0.289% |
| 40 | 13.782 | 1099 | 1102 | 1104 | rVV  | 3131574 | 3068196 | 25.37% | 3.951% |
| 41 | 13.885 | 1107 | 1111 | 1115 | rBV3 | 245784  | 569626  | 4.71%  | 0.734% |
| 42 | 14.011 | 1118 | 1122 | 1124 | rVB  | 487667  | 719591  | 5.95%  | 0.927% |
| 43 | 14.091 | 1126 | 1129 | 1130 | rBV  | 261742  | 334213  | 2.76%  | 0.430% |
| 44 | 14.137 | 1130 | 1133 | 1134 | rBV2 | 338490  | 404325  | 3.34%  | 0.521% |
| 45 | 14.240 | 1140 | 1142 | 1144 | rBV  | 223224  | 290025  | 2.40%  | 0.373% |
| 46 | 14.285 | 1144 | 1146 | 1147 | rVB  | 187545  | 257327  | 2.13%  | 0.331% |
| 47 | 14.605 | 1170 | 1174 | 1177 | rBV  | 259285  | 428598  | 3.54%  | 0.552% |
| 48 | 14.754 | 1183 | 1187 | 1188 | rBV2 | 210298  | 442128  | 3.66%  | 0.569% |
| 49 | 14.777 | 1188 | 1189 | 1191 | rVV  | 259372  | 320396  | 2.65%  | 0.413% |
| 50 | 14.811 | 1191 | 1192 | 1194 | rVV2 | 403630  | 561488  | 4.64%  | 0.723% |
| 51 | 14.857 | 1194 | 1196 | 1202 | rVB2 | 293670  | 499781  | 4.13%  | 0.644% |
| 52 | 15.040 | 1210 | 1212 | 1215 | rBV2 | 1509491 | 2834520 | 23.44% | 3.650% |
| 53 | 15.085 | 1215 | 1216 | 1221 | rVB  | 1207254 | 1494328 | 12.36% | 1.924% |
| 54 | 15.200 | 1224 | 1226 | 1228 | rBV  | 164979  | 278375  | 2.30%  | 0.358% |
| 55 | 15.325 | 1236 | 1237 | 1240 | rBV  | 132619  | 228269  | 1.89%  | 0.294% |
| 56 | 15.554 | 1254 | 1257 | 1258 | rBV  | 243319  | 376646  | 3.11%  | 0.485% |
| 57 | 15.588 | 1258 | 1260 | 1262 | rVB  | 162924  | 213371  | 1.76%  | 0.275% |
| 58 | 15.657 | 1262 | 1266 | 1267 | rBV4 | 112224  | 256619  | 2.12%  | 0.330% |
| 59 | 15.748 | 1272 | 1274 | 1276 | rVB  | 147684  | 195353  | 1.62%  | 0.252% |
| 60 | 15.806 | 1276 | 1279 | 1284 | rBV3 | 108707  | 193427  | 1.60%  | 0.249% |
| 61 | 16.434 | 1330 | 1334 | 1339 | rBV  | 1024297 | 2605908 | 21.55% | 3.356% |
| 62 | 16.583 | 1344 | 1347 | 1350 | rVV  | 168106  | 269763  | 2.23%  | 0.347% |
| 63 | 16.846 | 1367 | 1370 | 1374 | rVB  | 664274  | 1078277 | 8.92%  | 1.389% |
| 64 | 16.937 | 1374 | 1378 | 1381 | rBV  | 839029  | 1431551 | 11.84% | 1.844% |
| 65 | 17.017 | 1382 | 1385 | 1387 | rVV  | 579795  | 970249  | 8.02%  | 1.250% |
| 66 | 17.063 | 1387 | 1389 | 1393 | rVB2 | 260723  | 504479  | 4.17%  | 0.650% |
| 67 | 19.040 | 1558 | 1562 | 1570 | rVB2 | 364760  | 914175  | 7.56%  | 1.177% |
| 68 | 19.657 | 1612 | 1616 | 1623 | rVB  | 298549  | 762340  | 6.30%  | 0.982% |

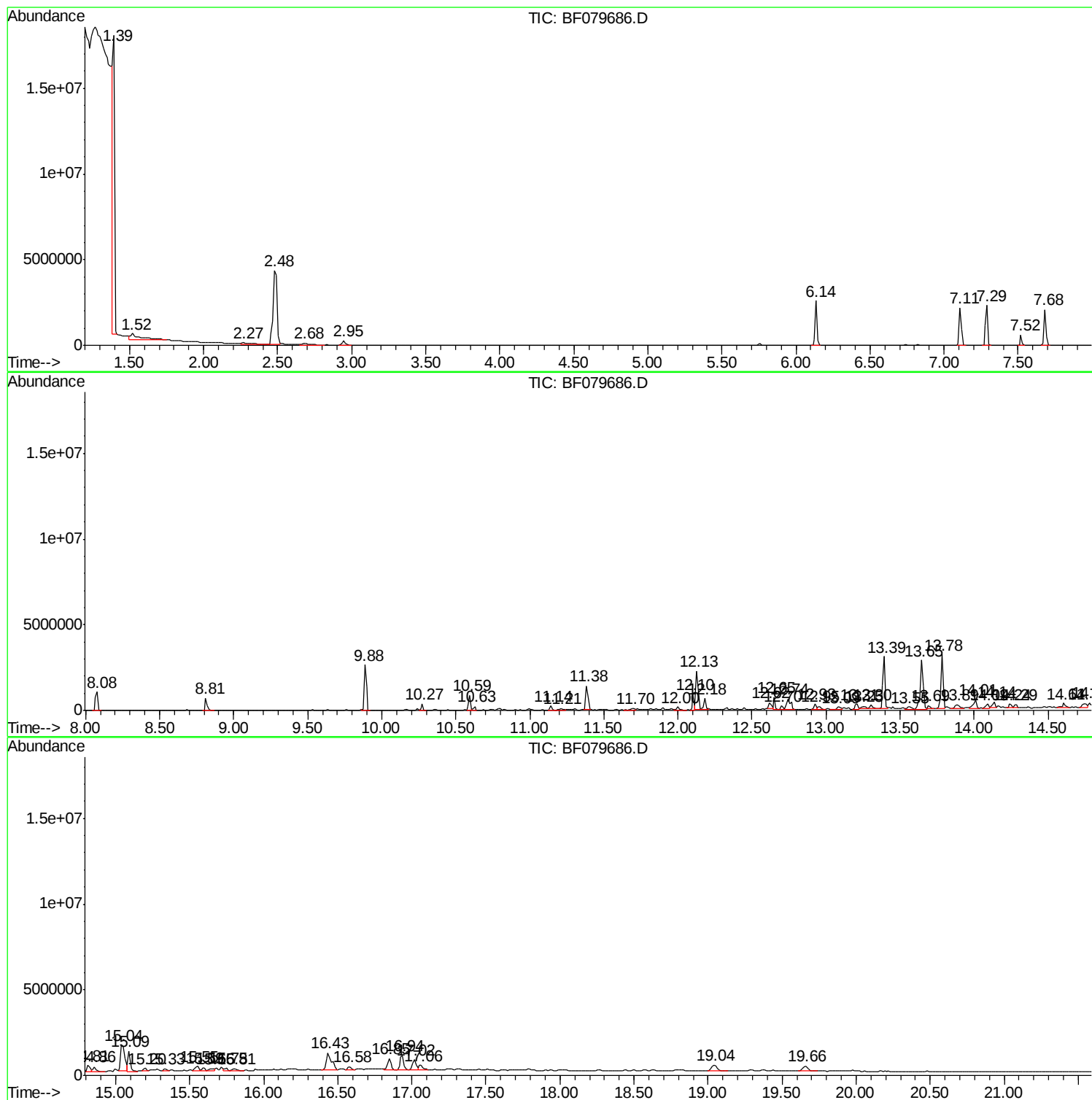
Sum of corrected areas: 77650709

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.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\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

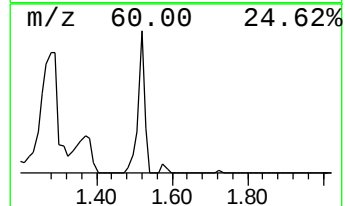
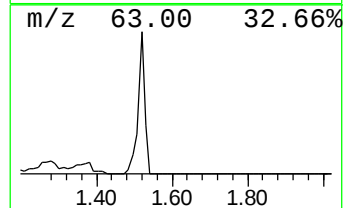
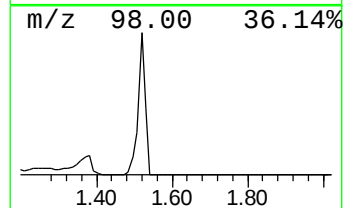
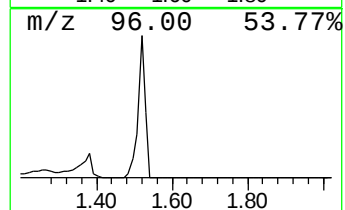
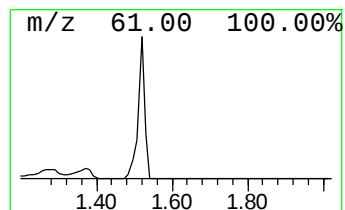
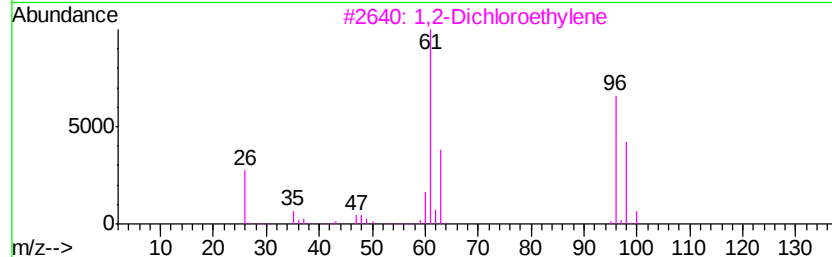
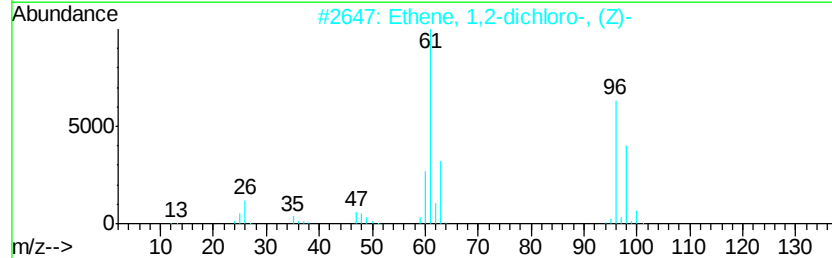
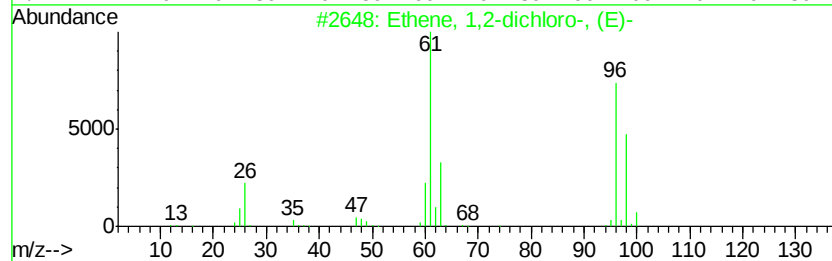
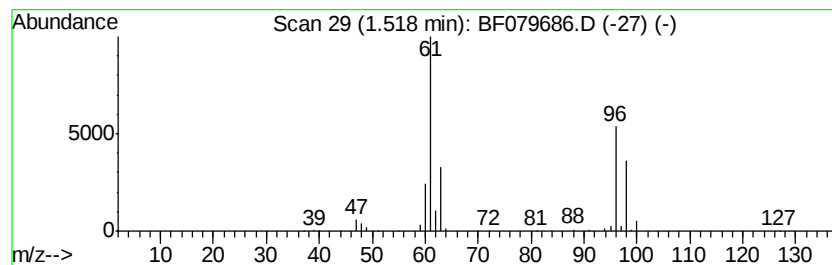
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Ethene, 1,2-dichloro-, (E)- Concentration Rank 5

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 1.52 | 65.25 ng | 1710540 | 1,4-Dichlorobenzene-d4 | 7.52 |

| Hit# | of | Tentative ID                | MW | MolForm | CAS#        | Qual |
|------|----|-----------------------------|----|---------|-------------|------|
| 1    | 5  | Ethene, 1,2-dichloro-, (E)- | 96 | C2H2Cl2 | 000156-60-5 | 94   |
| 2    |    | Ethene, 1,2-dichloro-, (Z)- | 96 | C2H2Cl2 | 000156-59-2 | 94   |
| 3    |    | 1,2-Dichloroethylene        | 96 | C2H2Cl2 | 000540-59-0 | 91   |
| 4    |    | Ethene, 1,1-dichloro-       | 96 | C2H2Cl2 | 000075-35-4 | 91   |
| 5    |    | Chloromethylmethyl sulfide  | 96 | C2H5ClS | 002373-51-5 | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

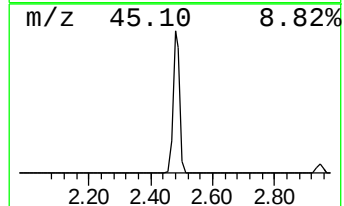
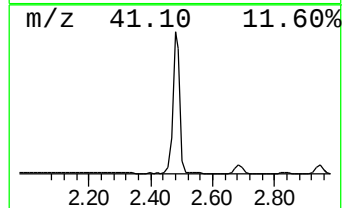
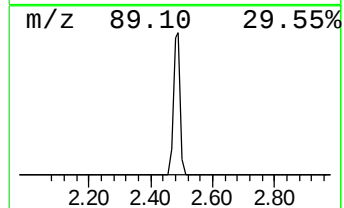
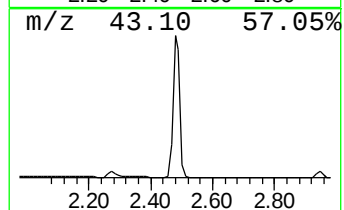
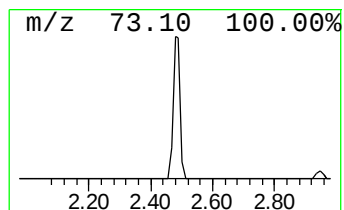
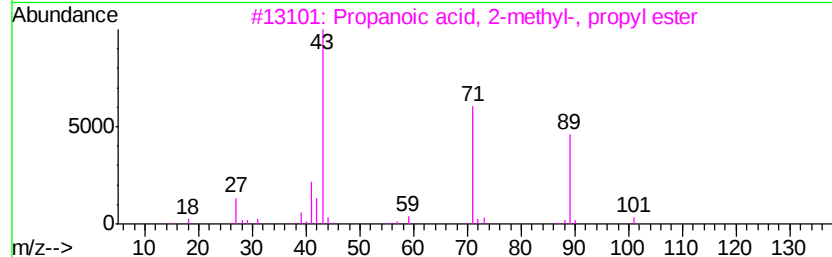
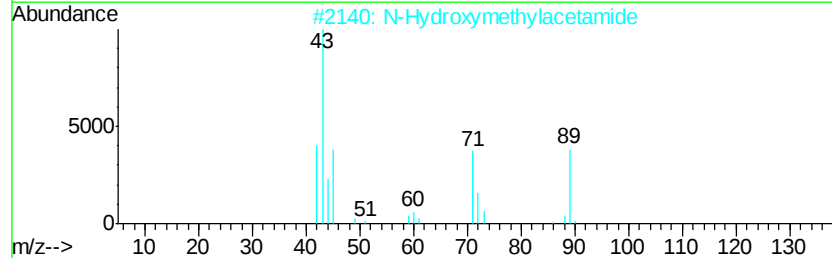
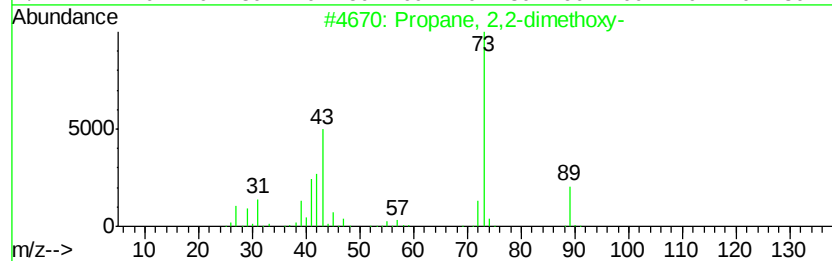
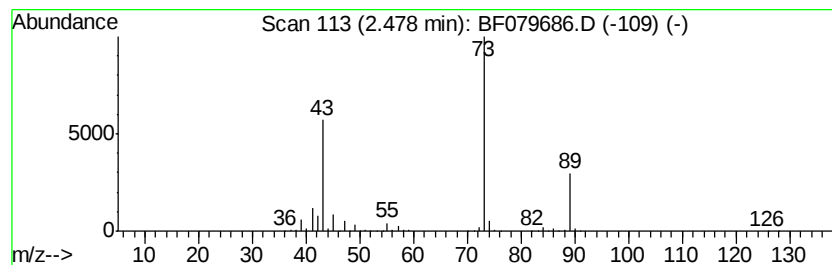
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Propane, 2,2-dimethoxy- Concentration Rank 2

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 2.48 | 282.02 ng | 7393520 | 1,4-Dichlorobenzene-d4 | 7.52 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Propane, 2,2-dimethoxy-              | 104 | C5H12O2 | 000077-76-9 | 59   |
| 2    |    | N-Hydroxymethylacetamide             | 89  | C3H7NO2 | 000625-51-4 | 17   |
| 3    |    | Propanoic acid, 2-methyl-, propyl... | 130 | C7H14O2 | 000644-49-5 | 12   |
| 4    |    | 2-Hydroxy-2-methylbutyric acid       | 118 | C5H10O3 | 003739-30-8 | 9    |
| 5    |    | Propane, 2-methoxy-2-methyl-         | 88  | C5H12O  | 001634-04-4 | 9    |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

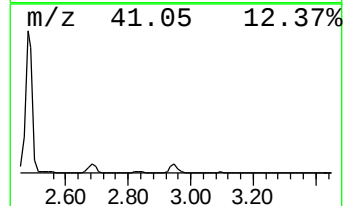
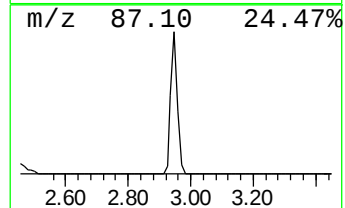
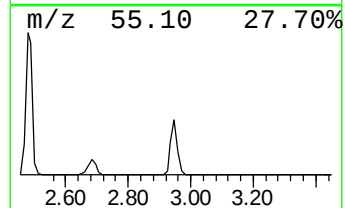
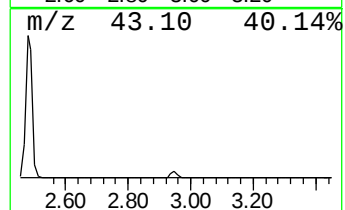
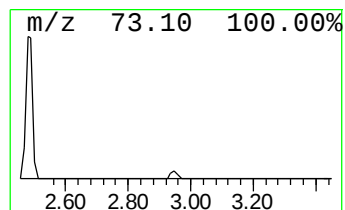
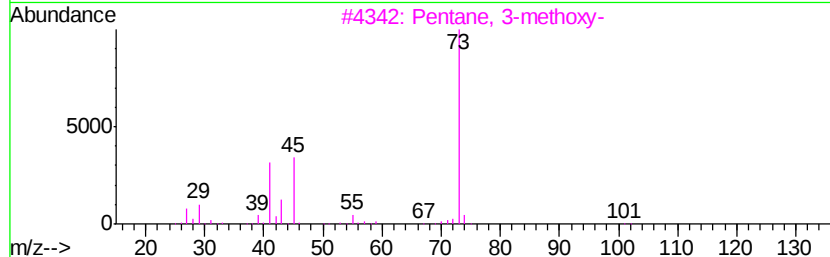
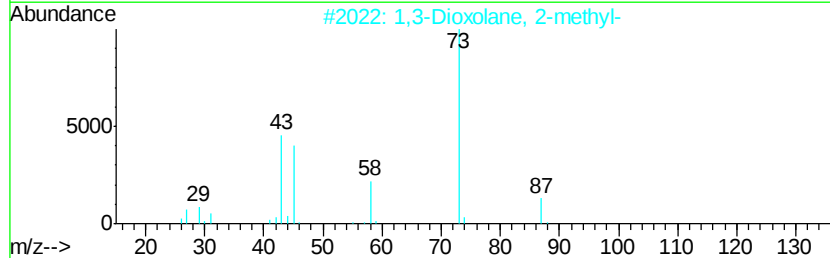
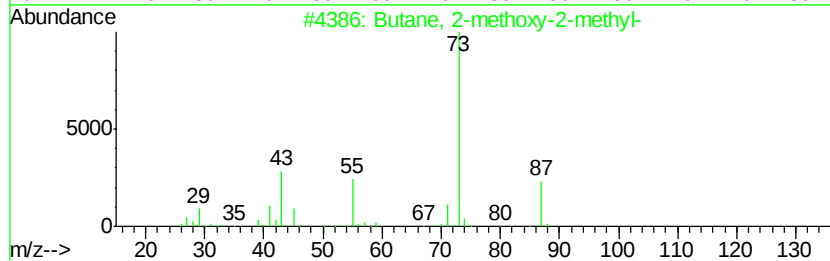
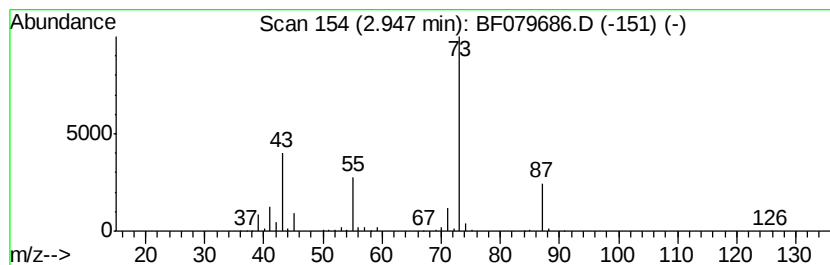
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Butane, 2-methoxy-2-methyl- Concentration Rank 8

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 2.95 | 14.42 ng | 378025 | 1,4-Dichlorobenzene-d4 | 7.52 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 37   |
| 3    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 12   |
| 4    |    | 1,4-Butanediamine           | 88  | C4H12N2 | 000110-60-1 | 9    |
| 5    |    | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

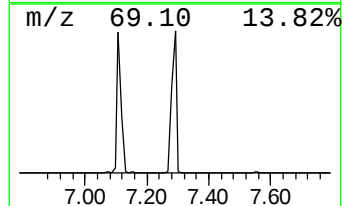
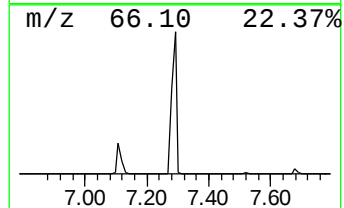
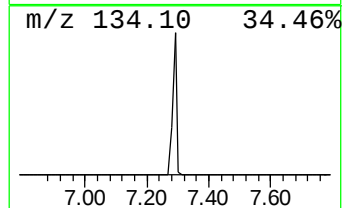
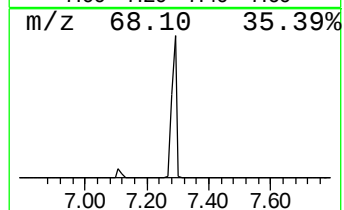
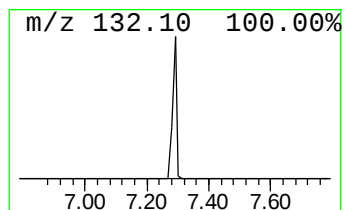
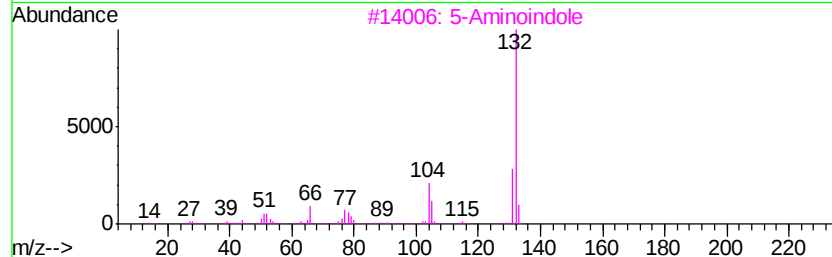
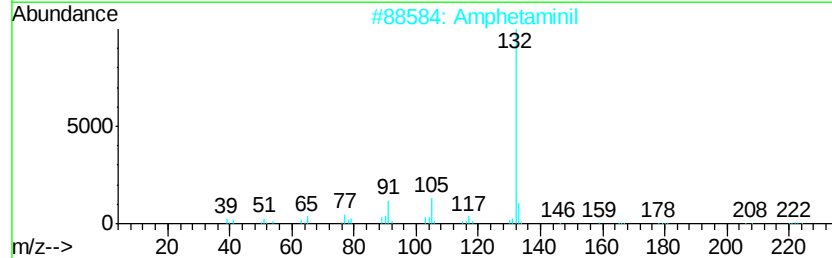
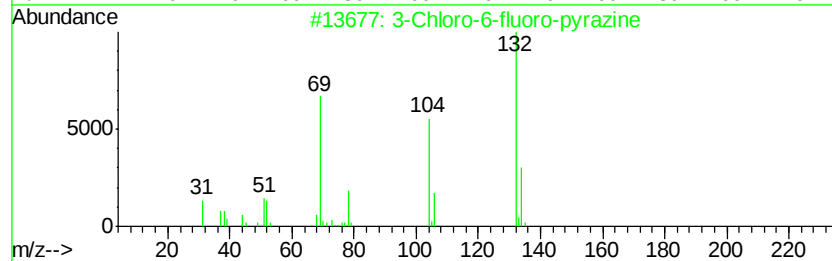
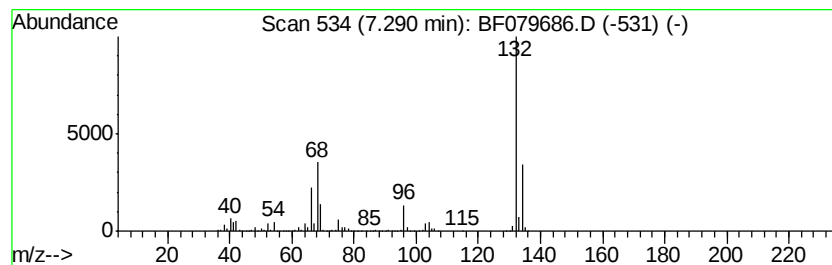
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 unknown7.29 Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.29 | 91.45 ng | 2397410 | 1,4-Dichlorobenzene-d4 | 7.52 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    |    | Amphetaminil                     | 250 | C17H18N2  | 017590-01-1  | 12   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 5    |    | Xenon                            | 132 | Xe        | 007440-63-3  | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

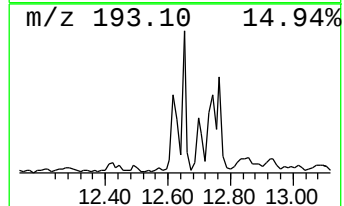
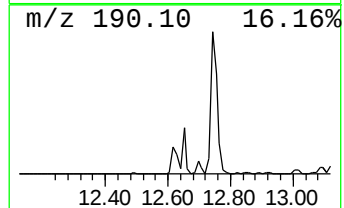
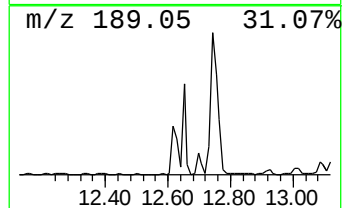
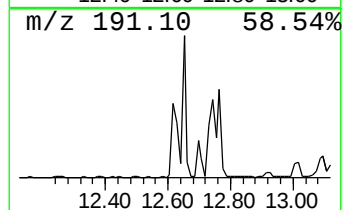
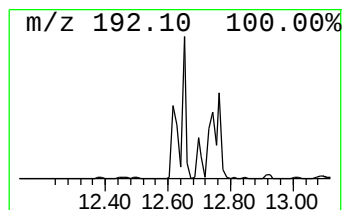
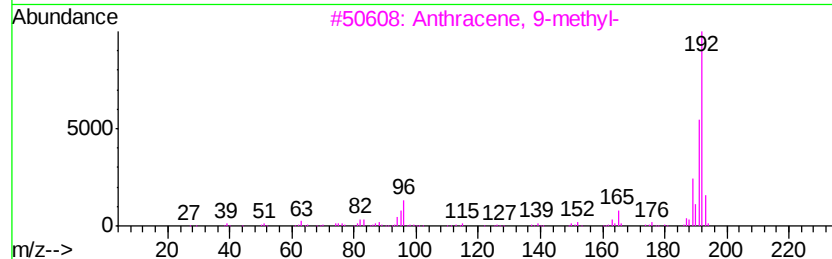
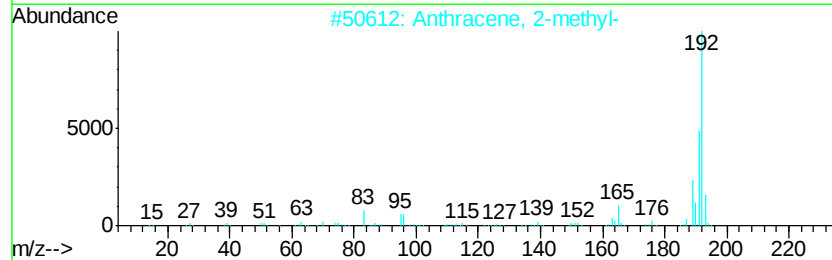
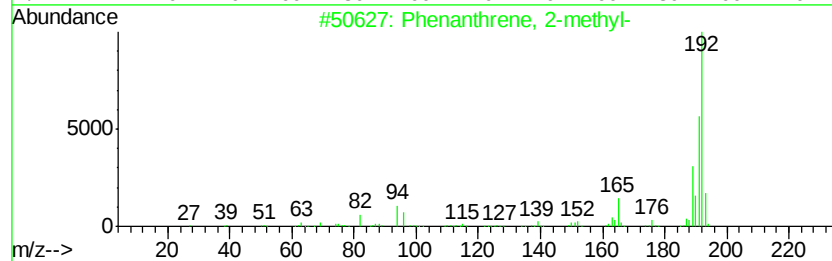
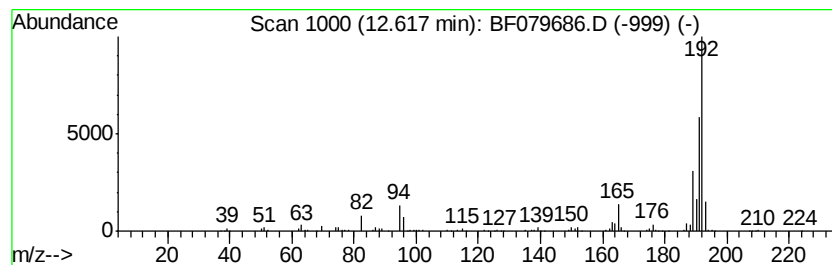
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 10

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.62 | 11.09 ng | 478297 | Phenanthrene-d10 | 12.10 |

| 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 | 95   |
| 3       |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |
| 4       |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 91   |
| 5       |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 90   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

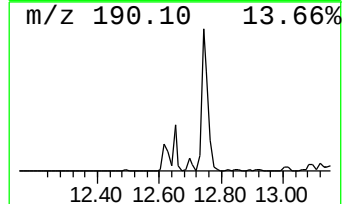
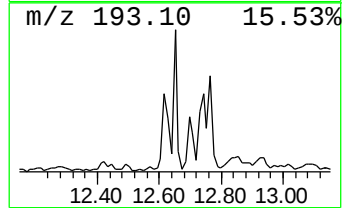
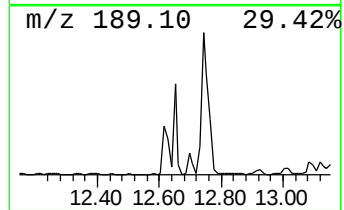
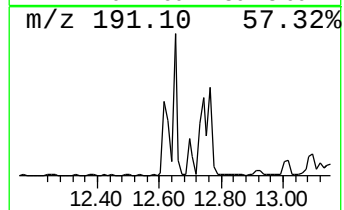
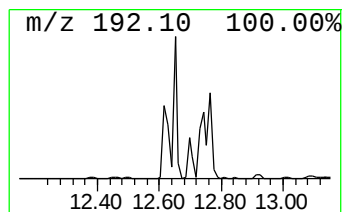
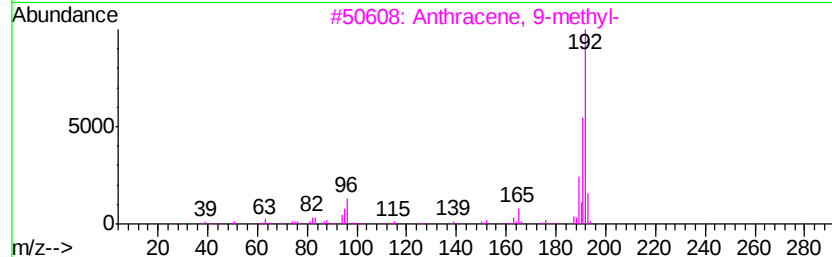
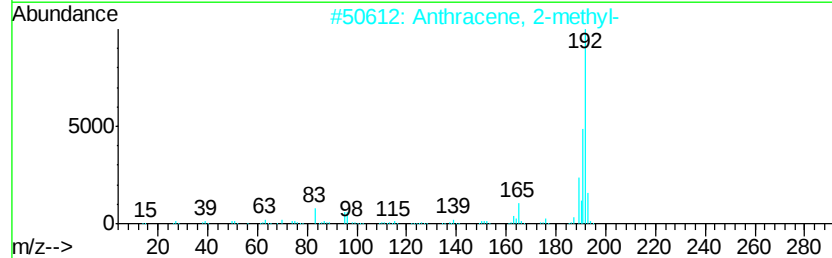
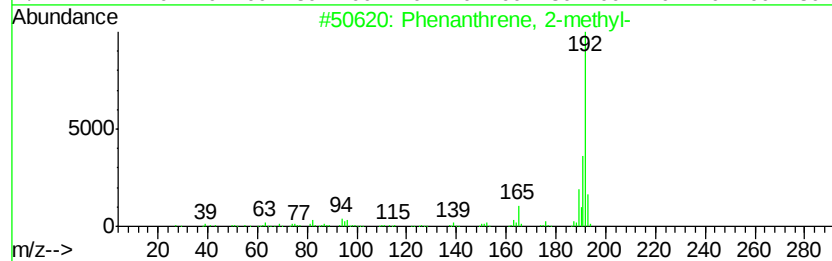
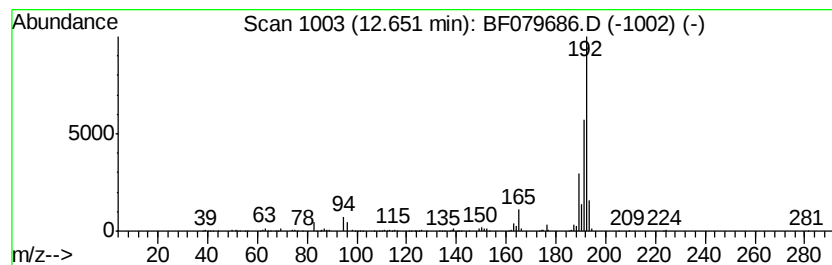
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Anthracene, 2-methyl- Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.65 | 12.43 ng | 536105 | Phenanthrene-d10 | 12.10 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2-methyl-               | 192 | C15H12  | 002531-84-2 | 96   |
| 2    |    | Anthracene, 2-methyl-                 | 192 | C15H12  | 000613-12-7 | 95   |
| 3    |    | Anthracene, 9-methyl-                 | 192 | C15H12  | 000779-02-2 | 95   |
| 4    |    | Phenanthrene, 3-methyl-               | 192 | C15H12  | 000832-71-3 | 94   |
| 5    |    | 1H-Cyclopropa[1]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 93   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

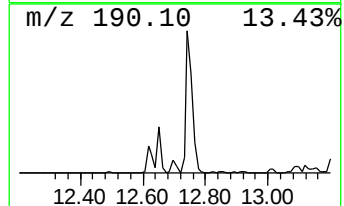
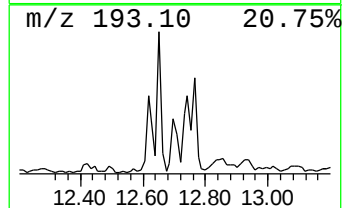
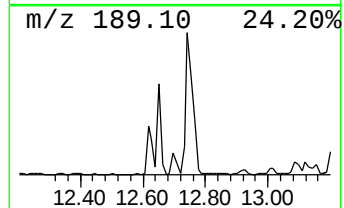
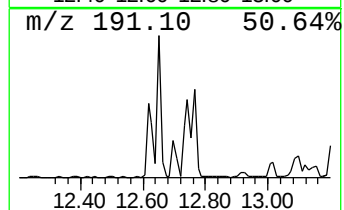
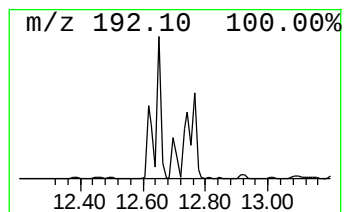
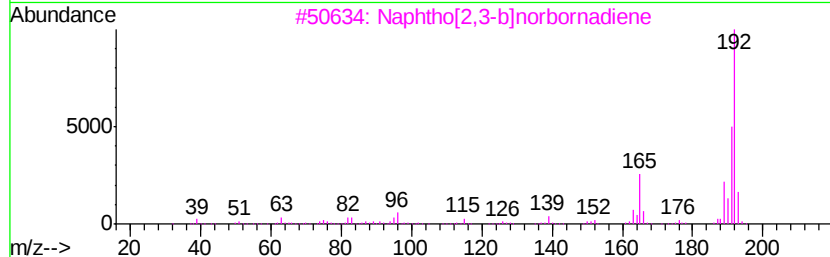
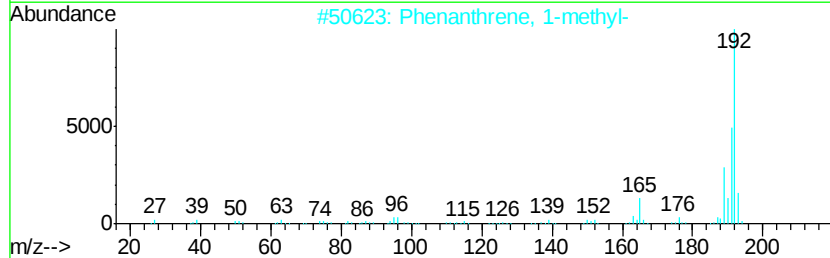
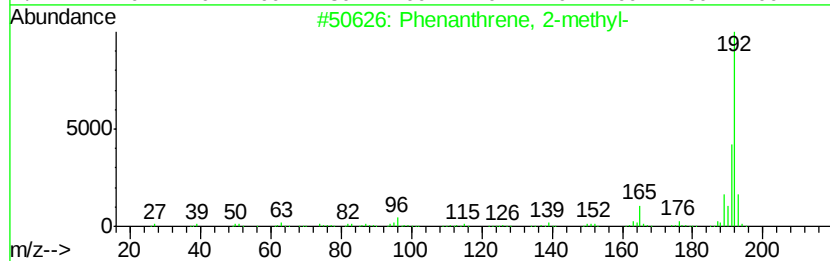
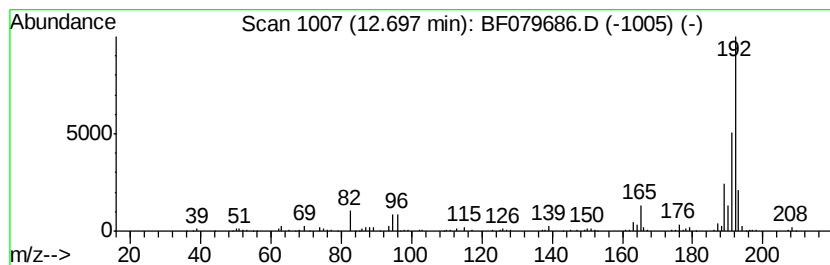
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Phenanthrene, 1-methyl- Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.70 | 5.60 ng | 241448 | Phenanthrene-d10 | 12.10 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl-     | 192 | C15H12  | 002531-84-2 | 96   |
| 2    |    |   | Phenanthrene, 1-methyl-     | 192 | C15H12  | 000832-69-9 | 93   |
| 3    |    |   | Naphtho[2,3-b]norbornadiene | 192 | C15H12  | 107426-38-0 | 91   |
| 4    |    |   | Phenanthrene, 3-methyl-     | 192 | C15H12  | 000832-71-3 | 91   |
| 5    |    |   | Anthracene, 1-methyl-       | 192 | C15H12  | 000610-48-0 | 91   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

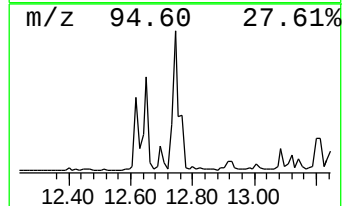
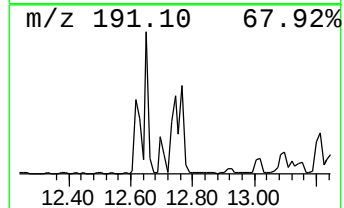
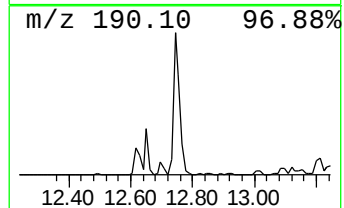
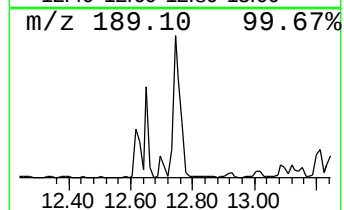
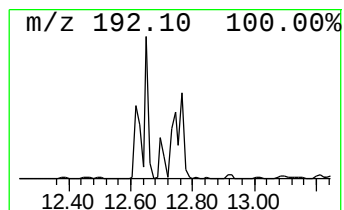
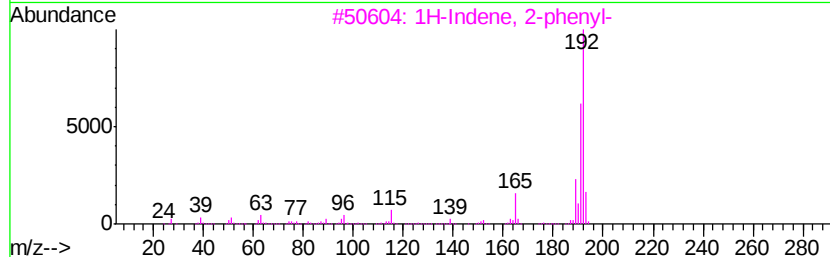
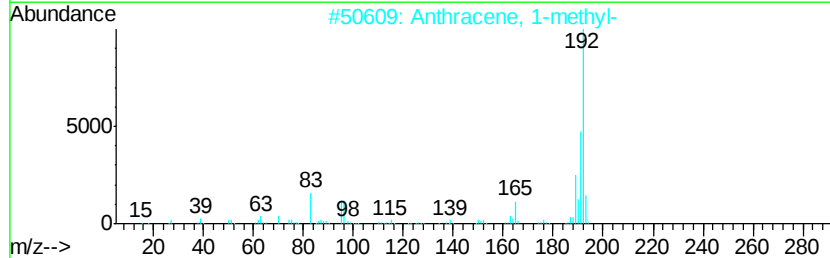
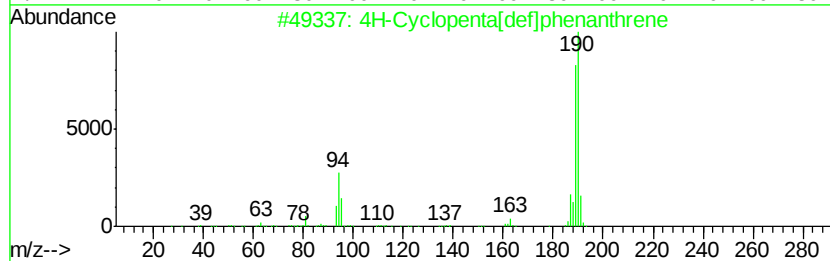
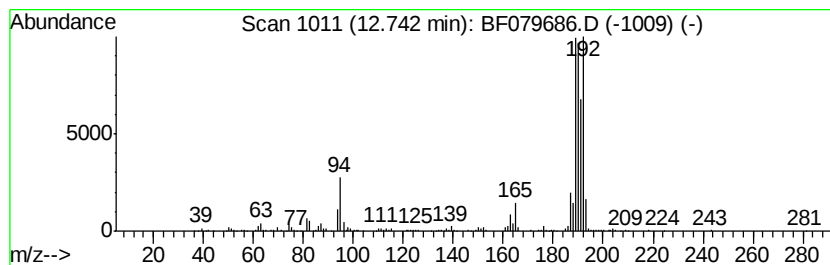
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.74 | 28.03 ng | 1209210 | Phenanthrene-d10 | 12.10 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10  | 000203-64-5 | 64   |
| 2    |    | Anthracene, 1-methyl-          | 192 | C15H12  | 000610-48-0 | 42   |
| 3    |    | 1H-Indene, 2-phenyl-           | 192 | C15H12  | 004505-48-0 | 42   |
| 4    |    | Anthracene, 2-methyl-          | 192 | C15H12  | 000613-12-7 | 42   |
| 5    |    | Phenanthrene, 1-methyl-        | 192 | C15H12  | 000832-69-9 | 42   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

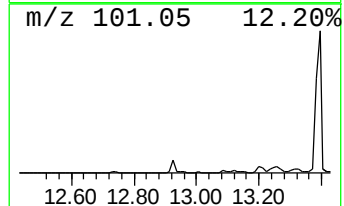
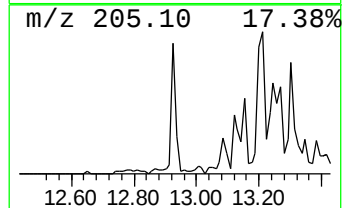
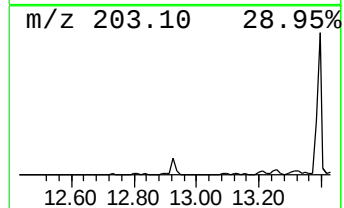
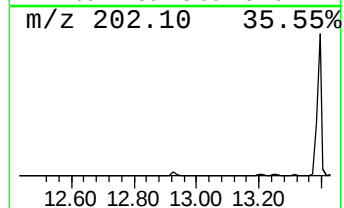
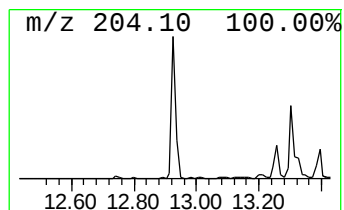
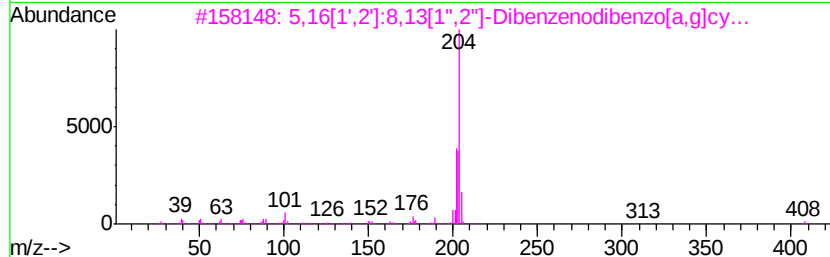
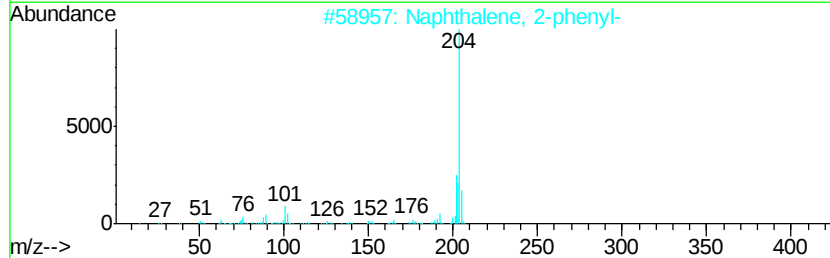
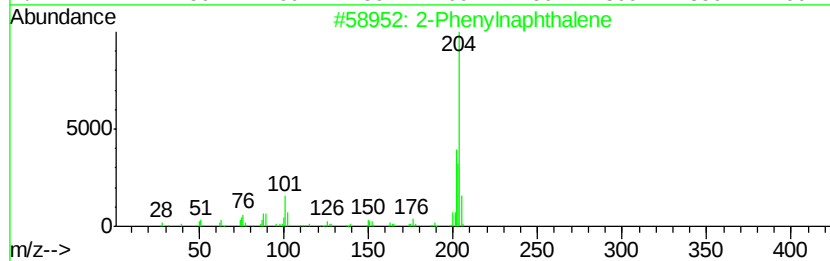
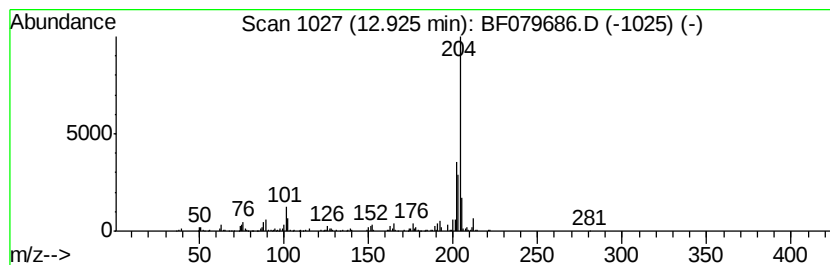
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 2-Phenylnaphthalene Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.93 | 7.10 ng | 306363 | Phenanthrene-d10 | 12.10 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | 2-Phenylnaphthalene                 | 204 | C16H12    | 035465-71-5 | 96   |
| 2    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12    | 000612-94-2 | 96   |
| 3    |    |   | 5,16[1',2']8,13[1'',2'']-Dibenz...  | 408 | C32H24    | 005672-97-9 | 87   |
| 4    |    |   | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12    | 006572-60-7 | 76   |
| 5    |    |   | 9,10-Bis(bromomethyl)anthracene     | 362 | C16H12Br2 | 034373-96-1 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

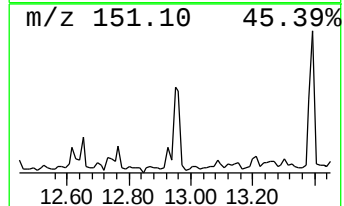
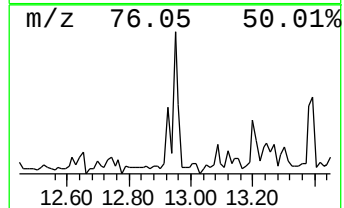
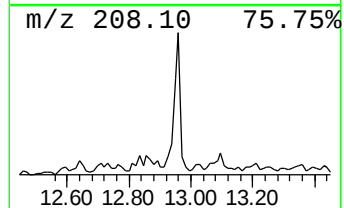
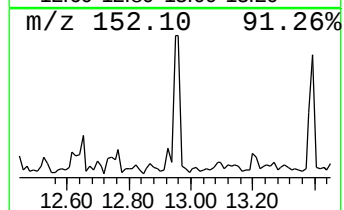
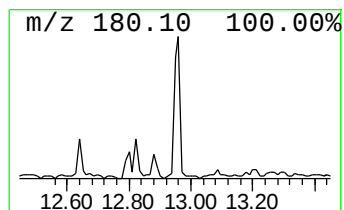
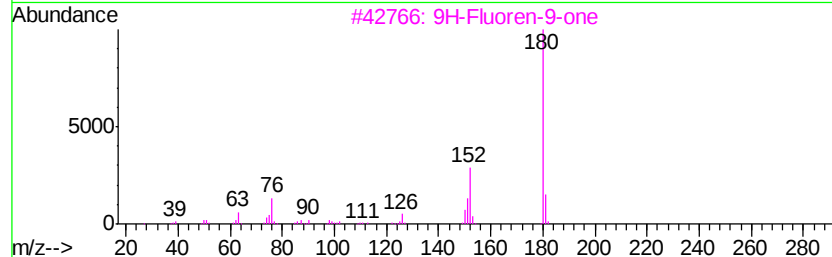
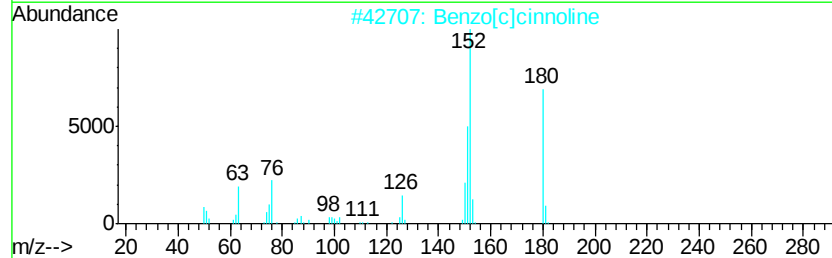
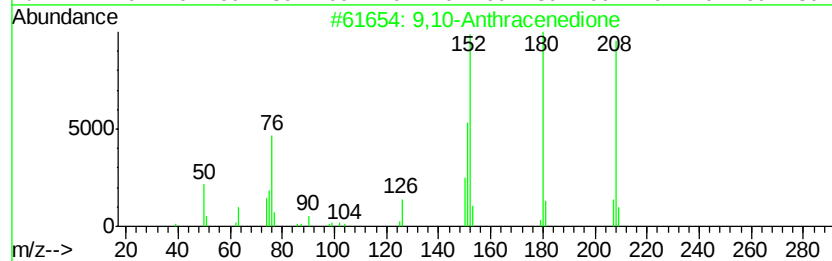
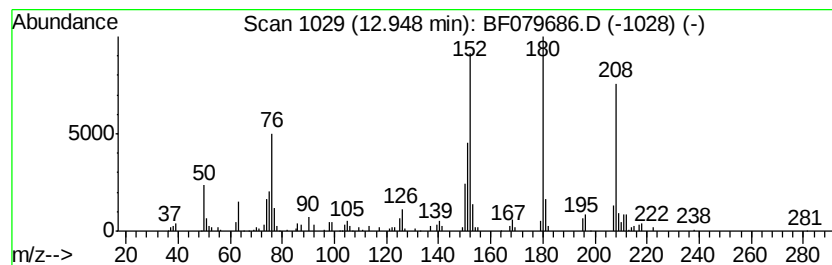
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 9,10-Anthracenedione Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.95 | 5.95 ng | 256843 | Phenanthrene-d10 | 12.10 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 99   |
| 2    |    |   | Benzo[c]cinnoline    | 180 | C12H8N2 | 000230-17-1 | 91   |
| 3    |    |   | 9H-Fluoren-9-one     | 180 | C13H8O  | 000486-25-9 | 78   |
| 4    |    |   | Benzo[h]cinnoline    | 180 | C12H8N2 | 000230-31-9 | 70   |
| 5    |    |   | 1H-Phenalen-1-one    | 180 | C13H8O  | 000548-39-0 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

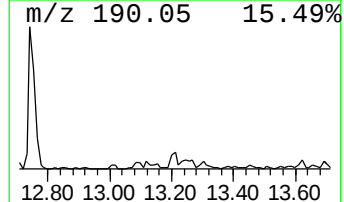
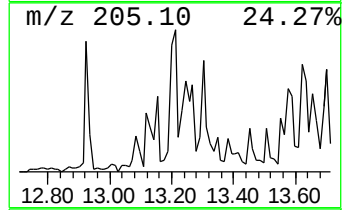
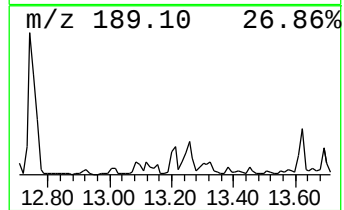
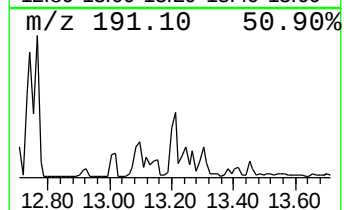
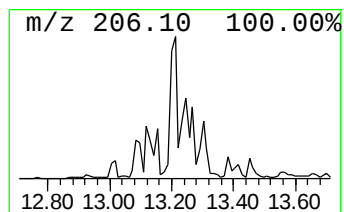
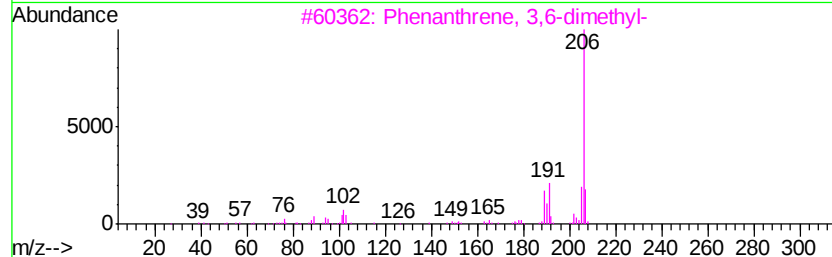
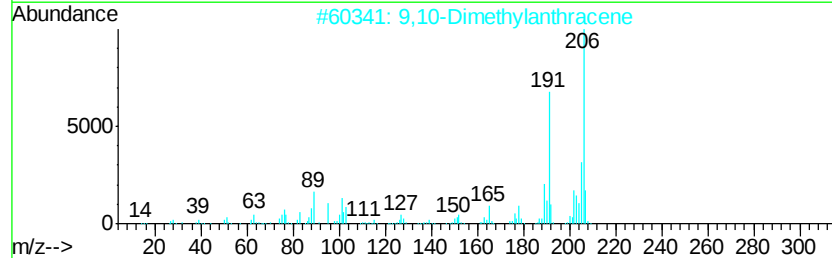
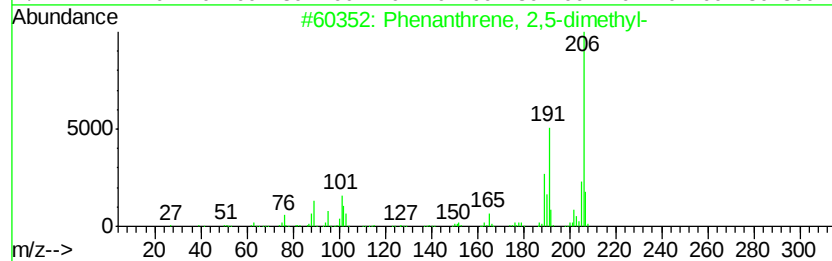
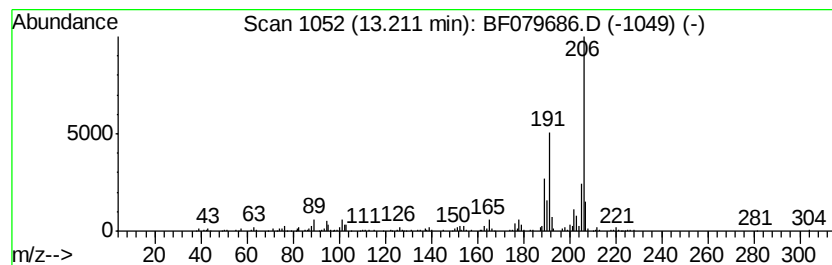
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Phenanthrene, 2,5-dimethyl- Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.21 | 9.98 ng | 430468 | Phenanthrene-d10 | 12.10 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|---------|---|-----------------------------|-----|---------|-------------|------|
| 1       |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 94   |
| 2       |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |
| 3       |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 91   |
| 4       |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 90   |
| 5       |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

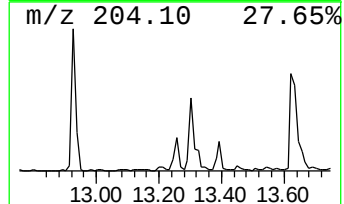
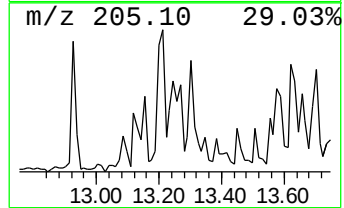
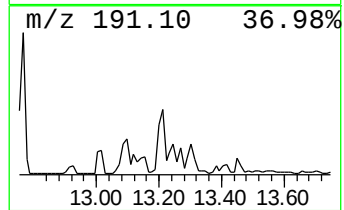
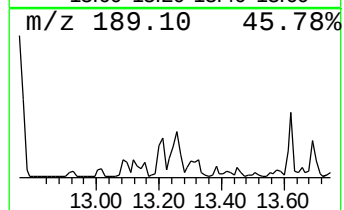
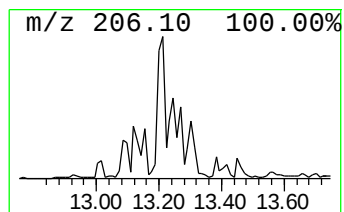
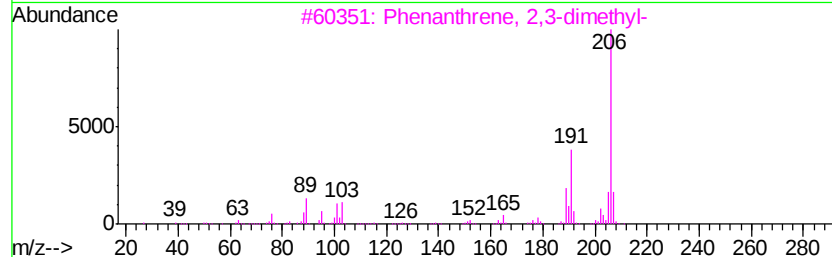
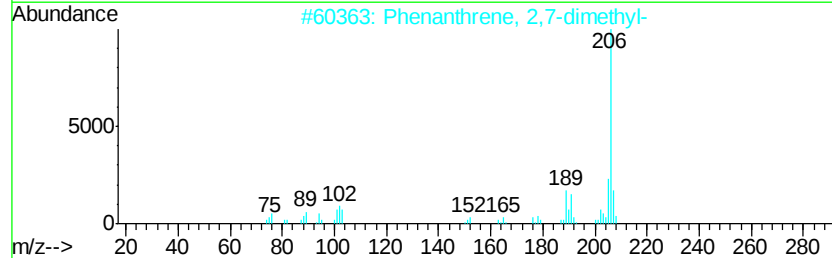
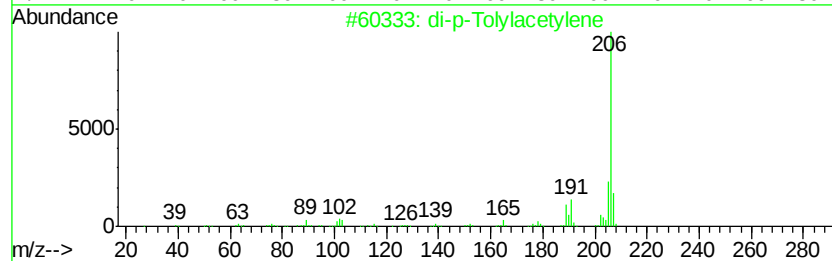
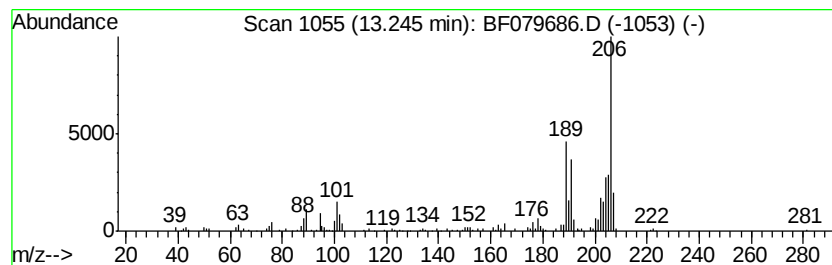
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 di-p-Tolylacetylene Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.25 | 8.25 ng | 356094 | Phenanthrene-d10 | 12.10 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | di-p-Tolylacetylene                | 206 | C16H14  | 002789-88-0 | 93   |
| 2    |    | Phenanthrene, 2,7-dimethyl-        | 206 | C16H14  | 001576-69-8 | 87   |
| 3    |    | Phenanthrene, 2,3-dimethyl-        | 206 | C16H14  | 003674-65-5 | 86   |
| 4    |    | 9,10-Dimethylantracene             | 206 | C16H14  | 000781-43-1 | 76   |
| 5    |    | Naphthalene, 1,2-dihydro-4-phenyl- | 206 | C16H14  | 007469-40-1 | 70   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

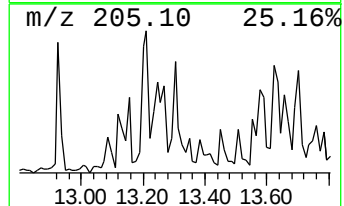
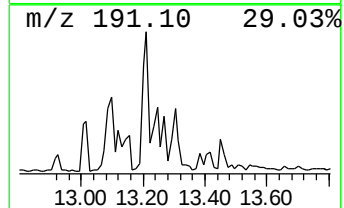
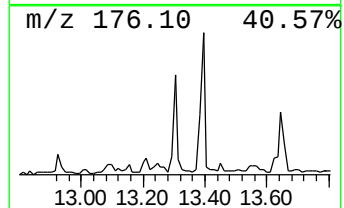
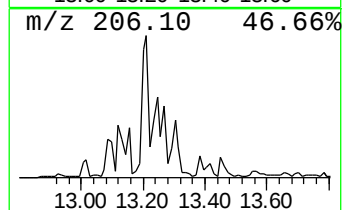
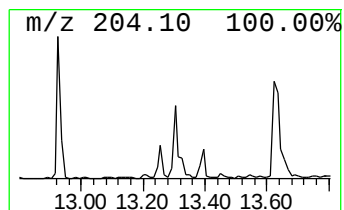
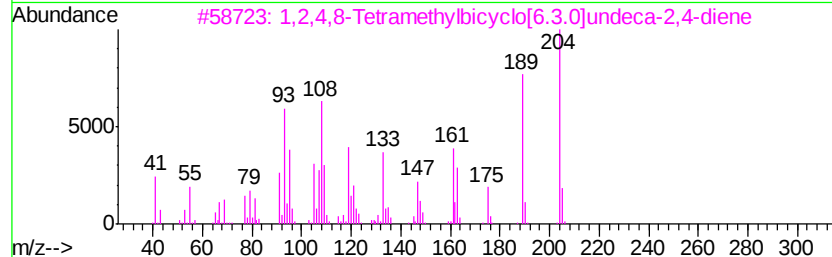
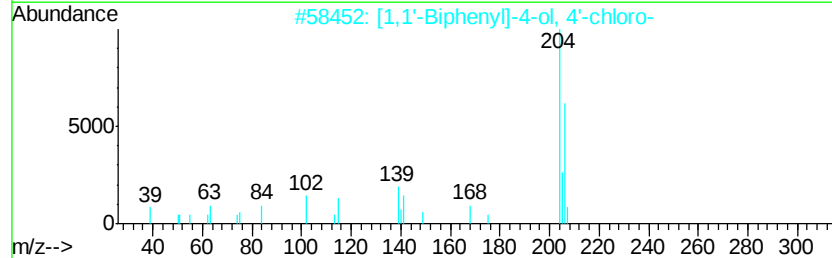
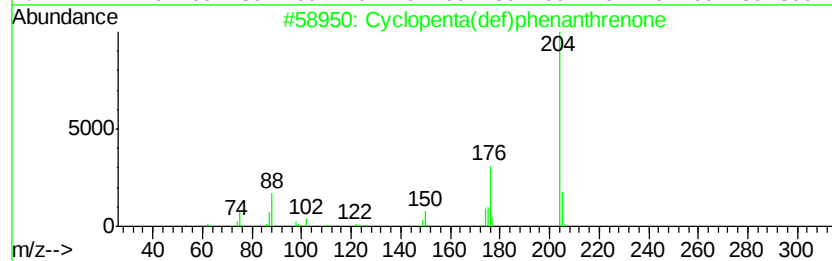
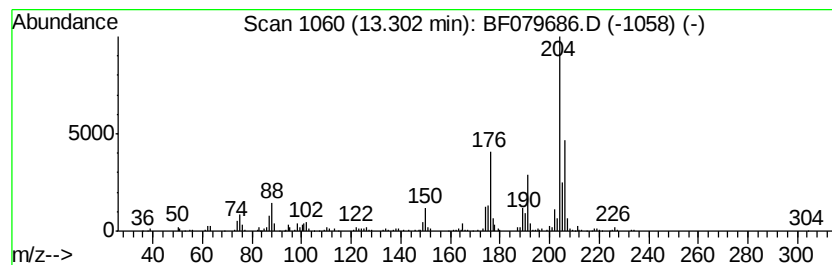
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Cyclopenta(def)phenanthrenone Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.30 | 8.47 ng | 365644 | Phenanthrene-d10 | 12.10 |

| Hit# | of | Tentative ID                                      | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrenone                     | 204 | C15H8O   | 005737-13-3 | 90   |
| 2    |    | [1,1'-Biphenyl]-4-ol, 4'-chloro-                  | 204 | C12H9ClO | 028034-99-3 | 49   |
| 3    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene | 204 | C15H24   | 137235-51-9 | 41   |
| 4    |    | [1,1'-Biphenyl]-4-ol, 3-chloro-                   | 204 | C12H9ClO | 000092-04-6 | 38   |
| 5    |    | Benzene, 1-chloro-3-phenoxy-                      | 204 | C12H9ClO | 006452-49-9 | 37   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

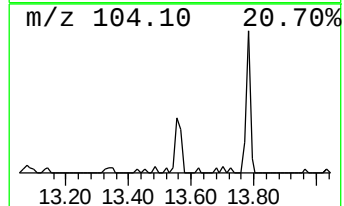
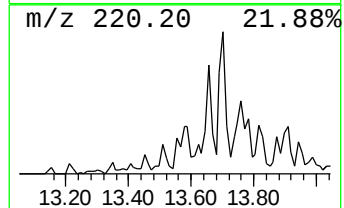
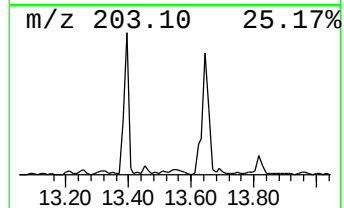
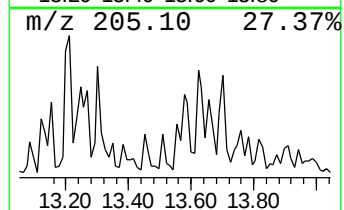
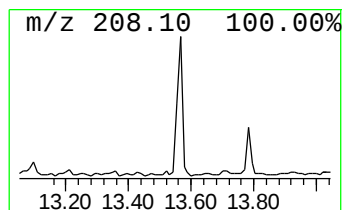
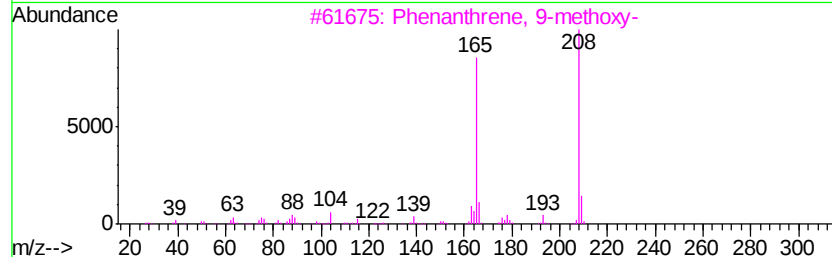
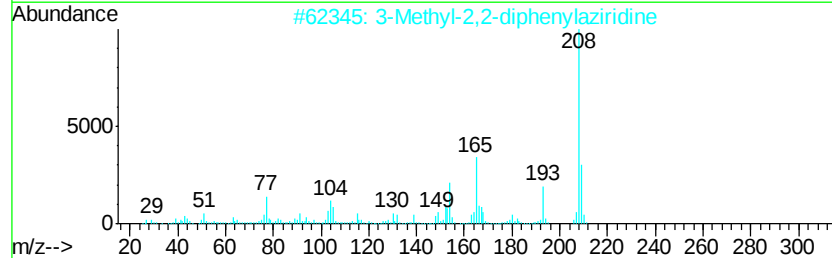
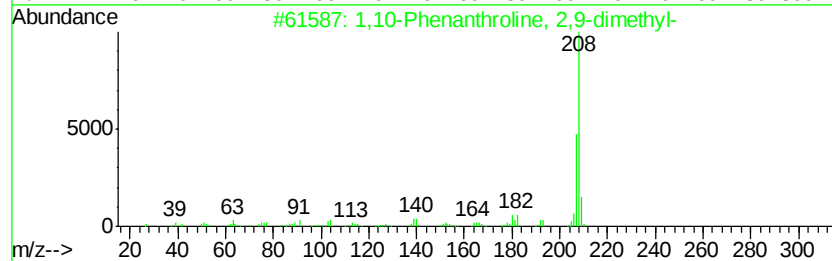
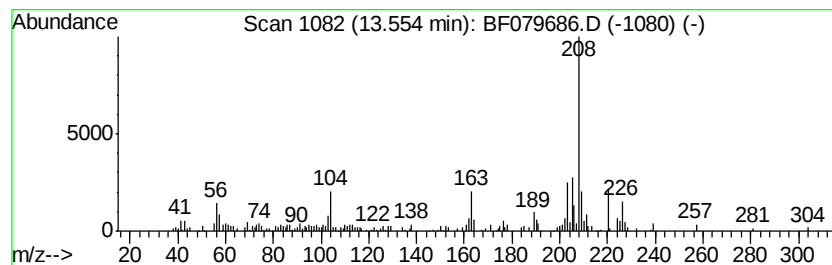
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 unknown13.55 Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.55 | 5.96 ng | 256938 | Phenanthrene-d10 | 12.10 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1,10-Phenanthroline, 2,9-dimethyl-  | 208 | C14H12N2  | 000484-11-7  | 38   |
| 2    |    |   | 3-Methyl-2,2-diphenylaziridine      | 209 | C15H15N   | 007764-13-8  | 35   |
| 3    |    |   | Phenanthrene, 9-methoxy-            | 208 | C15H12O   | 005085-74-5  | 35   |
| 4    |    |   | 4,7-Dimethyl-1,10-phenanthroline    | 208 | C14H12N2  | 003248-05-3  | 32   |
| 5    |    |   | 2-p-Tolyl-2,3-dihydro-1H-benzo[1... | 208 | C13H13BN2 | 1000296-11-8 | 32   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
Data File : BF079686.D  
Acq On : 4 Jun 2015 4:28  
Operator : TP/IZ  
Sample : G2408-25  
Misc :  
ALS Vial : 32 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

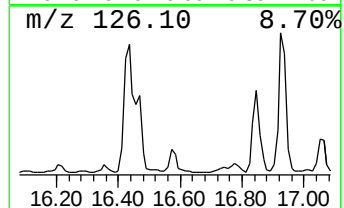
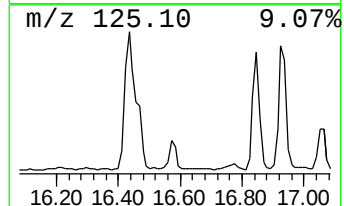
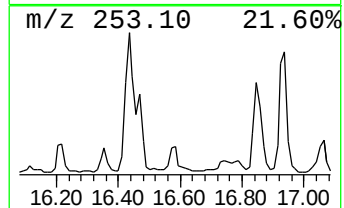
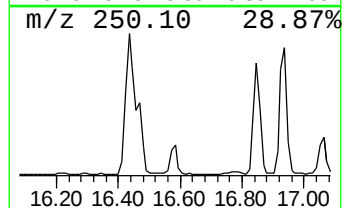
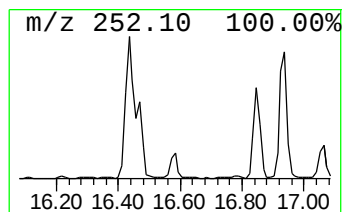
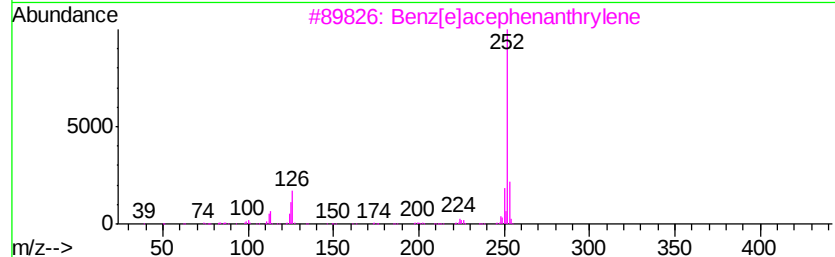
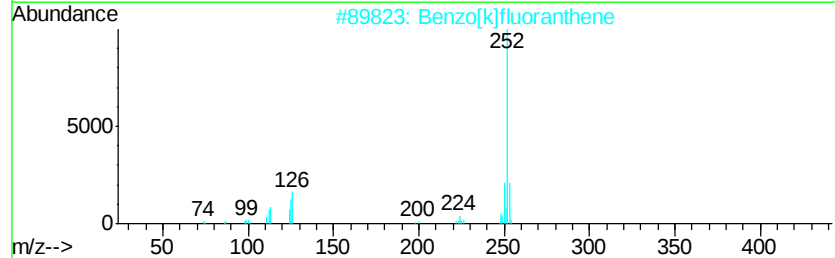
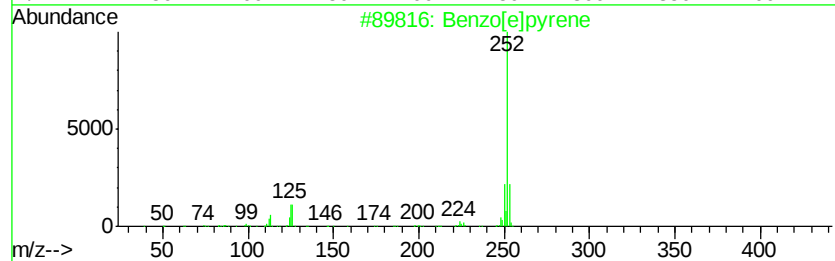
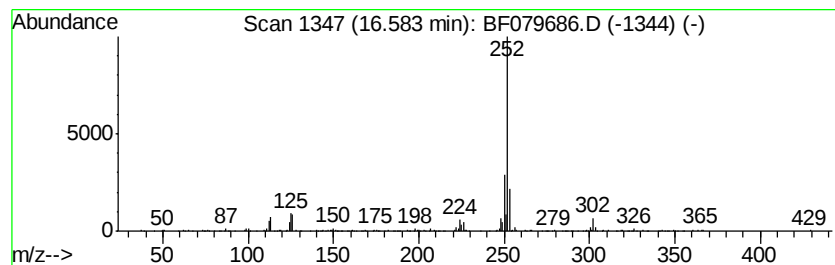
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Benzo[e]pyrene Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.58 | 5.56 ng | 269763 | Perylene-d12     | 17.02 |

| Hit# of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|---------|---|---------------------------|-----|---------|-------------|------|
| 1       |   | Benzo[e]pyrene            | 252 | C20H12  | 000192-97-2 | 95   |
| 2       |   | Benzo[k]fluoranthene      | 252 | C20H12  | 000207-08-9 | 94   |
| 3       |   | Benzo[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 90   |
| 4       |   | Benzo[a]pyrene            | 252 | C20H12  | 000050-32-8 | 90   |
| 5       |   | Benzo[j]fluoranthene      | 252 | C20H12  | 000205-82-3 | 89   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060315\  
 Data File : BF079686.D  
 Acq On : 4 Jun 2015 4:28  
 Operator : TP/IZ  
 Sample : G2408-25  
 Misc :  
 ALS Vial : 32 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-14D

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060315.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 |
| Ethene, 1,2-dichl... | 1.52  | 65.3    | ng    | 1710540  | 1                     | 7.52  | 524321 | 20.0 |
| Propane, 2,2-dime... | 2.48  | 282.0   | ng    | 7393520  | 1                     | 7.52  | 524321 | 20.0 |
| Butane, 2-methoxy... | 2.95  | 14.4    | ng    | 378025   | 1                     | 7.52  | 524321 | 20.0 |
| unknown7.29          | 7.29  | 91.5    | ng    | 2397410  | 1                     | 7.52  | 524321 | 20.0 |
| Phenanthrene, 2-m... | 12.62 | 11.1    | ng    | 478297   | 4                     | 12.10 | 862932 | 20.0 |
| Anthracene, 2-met... | 12.65 | 12.4    | ng    | 536105   | 4                     | 12.10 | 862932 | 20.0 |
| Phenanthrene, 1-m... | 12.70 | 5.6     | ng    | 241448   | 4                     | 12.10 | 862932 | 20.0 |
| 4H-Cyclopenta[def... | 12.74 | 28.0    | ng    | 1209210  | 4                     | 12.10 | 862932 | 20.0 |
| 2-Phenyl naphthalene | 12.93 | 7.1     | ng    | 306363   | 4                     | 12.10 | 862932 | 20.0 |
| 9,10-Anthracenedione | 12.95 | 6.0     | ng    | 256843   | 4                     | 12.10 | 862932 | 20.0 |
| Phenanthrene, 2,5... | 13.21 | 10.0    | ng    | 430468   | 4                     | 12.10 | 862932 | 20.0 |
| di-p-Tolylacetylene  | 13.25 | 8.3     | ng    | 356094   | 4                     | 12.10 | 862932 | 20.0 |
| Cyclopenta(def)ph... | 13.30 | 8.5     | ng    | 365644   | 4                     | 12.10 | 862932 | 20.0 |
| unknown13.55         | 13.55 | 6.0     | ng    | 256938   | 4                     | 12.10 | 862932 | 20.0 |
| Benzo[e]pyrene       | 16.58 | 5.6     | ng    | 269763   | 6                     | 17.02 | 970249 | 20.0 |