

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012215\  
 Data File : BF077032.D  
 Acq On : 23 Jan 2015 5:08  
 Operator : TP/IZ  
 Sample : G1148-03  
 Misc : S  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 FK-GF-02-004-B

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-BF011415.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         | 1.350       | 21            | 23          | 28           | rBV2     | 6163           | 11702         | 2.67%           | 0.232%        |
| 2         | 1.430       | 28            | 30          | 41           | rVB      | 58310          | 134774        | 30.74%          | 2.667%        |
| 3         | 2.733       | 141           | 144         | 150          | rBV      | 15590          | 37339         | 8.52%           | 0.739%        |
| 4         | 3.944       | 247           | 250         | 257          | rBV      | 66896          | 148862        | 33.96%          | 2.946%        |
| 5         | 4.893       | 330           | 333         | 340          | rBV      | 200384         | 380996        | 86.91%          | 7.541%        |
| 6         | 7.248       | 536           | 539         | 544          | rBV      | 233552         | 388457        | 88.61%          | 7.688%        |
| 7         | 7.339       | 544           | 547         | 553          | rVB      | 254697         | 424939        | 96.93%          | 8.410%        |
| 8         | 7.853       | 588           | 592         | 596          | rVB      | 90218          | 137948        | 31.47%          | 2.730%        |
| 9         | 8.173       | 617           | 620         | 624          | rBV      | 193620         | 304000        | 69.34%          | 6.017%        |
| 10        | 9.145       | 702           | 705         | 713          | rBV      | 154899         | 253656        | 57.86%          | 5.020%        |
| 11        | 10.756      | 843           | 846         | 850          | rBV      | 122218         | 191384        | 43.65%          | 3.788%        |
| 12        | 12.231      | 973           | 975         | 979          | rVB      | 12323          | 18514         | 4.22%           | 0.366%        |
| 13        | 13.408      | 1075          | 1078        | 1082         | rBV      | 287789         | 430694        | 98.24%          | 8.524%        |
| 14        | 14.483      | 1169          | 1172        | 1176         | rBV      | 21934          | 35772         | 8.16%           | 0.708%        |
| 15        | 14.883      | 1204          | 1207        | 1211         | rVB      | 132699         | 205318        | 46.83%          | 4.064%        |
| 16        | 16.140      | 1315          | 1317        | 1320         | rBV      | 5357           | 7614          | 1.74%           | 0.151%        |
| 17        | 16.528      | 1348          | 1351        | 1353         | rBV      | 69150          | 80588         | 18.38%          | 1.595%        |
| 18        | 16.586      | 1353          | 1356        | 1359         | rVB      | 190572         | 269496        | 61.47%          | 5.334%        |
| 19        | 17.694      | 1450          | 1453        | 1455         | rBV      | 158425         | 215841        | 49.23%          | 4.272%        |
| 20        | 17.729      | 1455          | 1456        | 1458         | rVB      | 15112          | 11343         | 2.59%           | 0.225%        |
| 21        | 19.374      | 1598          | 1600        | 1602         | rVB      | 24885          | 31798         | 7.25%           | 0.629%        |
| 22        | 19.649      | 1622          | 1624        | 1627         | rVB      | 24600          | 30981         | 7.07%           | 0.613%        |
| 23        | 19.820      | 1637          | 1639        | 1642         | rBV      | 11593          | 17789         | 4.06%           | 0.352%        |
| 24        | 19.923      | 1646          | 1648        | 1651         | rVB      | 434486         | 438404        | 100.00%         | 8.677%        |
| 25        | 21.192      | 1755          | 1759        | 1761         | rBV      | 222310         | 333611        | 76.10%          | 6.603%        |
| 26        | 21.878      | 1815          | 1819        | 1820         | rBV2     | 3547           | 5960          | 1.36%           | 0.118%        |
| 27        | 21.969      | 1824          | 1827        | 1835         | rVB3     | 5603           | 17937         | 4.09%           | 0.355%        |
| 28        | 22.335      | 1856          | 1859        | 1863         | rBV2     | 3273           | 5898          | 1.35%           | 0.117%        |
| 29        | 22.426      | 1865          | 1867        | 1871         | rVB      | 16682          | 28581         | 6.52%           | 0.566%        |
| 30        | 22.723      | 1888          | 1893        | 1895         | rBV2     | 9961           | 21551         | 4.92%           | 0.427%        |
| 31        | 22.781      | 1895          | 1898        | 1900         | rVV      | 12164          | 16349         | 3.73%           | 0.324%        |
| 32        | 22.838      | 1900          | 1903        | 1908         | rBV      | 181685         | 212605        | 48.50%          | 4.208%        |
| 33        | 23.066      | 1920          | 1923        | 1925         | rBV2     | 2841           | 6076          | 1.39%           | 0.120%        |
| 34        | 23.363      | 1943          | 1949        | 1951         | rBV7     | 2717           | 9154          | 2.09%           | 0.181%        |

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ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
FK-GF-02-004-B

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-BF011415.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |       |       |       |        |
|----|--------|------|------|------|------|-------|-------|-------|--------|
| 35 | 23.421 | 1951 | 1954 | 1956 | rBV3 | 10381 | 15197 | 3.47% | 0.301% |
| 36 | 23.466 | 1956 | 1958 | 1964 | rVB5 | 12522 | 30519 | 6.96% | 0.604% |
| 37 | 23.798 | 1983 | 1987 | 1989 | rBV3 | 6175  | 11268 | 2.57% | 0.223% |
| 38 | 23.935 | 1997 | 1999 | 2005 | rVB3 | 8414  | 16411 | 3.74% | 0.325% |
| 39 | 24.164 | 2017 | 2019 | 2021 | rBV3 | 4777  | 7710  | 1.76% | 0.153% |
| 40 | 24.221 | 2021 | 2024 | 2027 | rVB4 | 9325  | 18980 | 4.33% | 0.376% |
| 41 | 24.381 | 2035 | 2038 | 2041 | rBV5 | 6162  | 14064 | 3.21% | 0.278% |
| 42 | 24.552 | 2049 | 2053 | 2057 | rBV5 | 15095 | 36275 | 8.27% | 0.718% |
| 43 | 24.701 | 2064 | 2066 | 2069 | rVB4 | 3493  | 4933  | 1.13% | 0.098% |
| 44 | 24.872 | 2077 | 2081 | 2083 | rBV5 | 5667  | 12970 | 2.96% | 0.257% |
| 45 | 25.009 | 2091 | 2093 | 2096 | rBV3 | 3597  | 8029  | 1.83% | 0.159% |
| 46 | 25.638 | 2146 | 2148 | 2151 | rVB3 | 4513  | 5821  | 1.33% | 0.115% |
| 47 | 25.855 | 2165 | 2167 | 2170 | rBV4 | 2115  | 4423  | 1.01% | 0.088% |

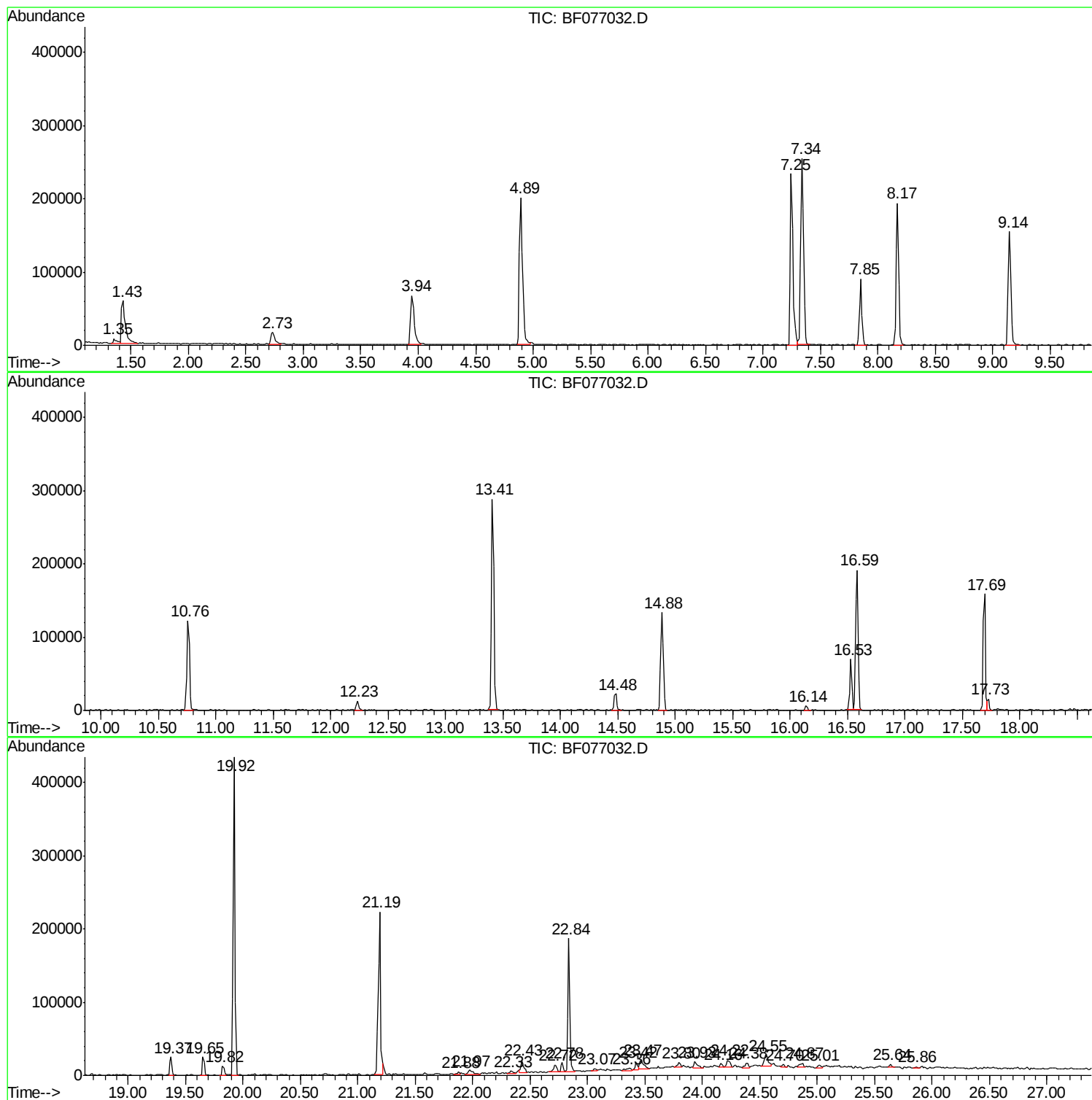
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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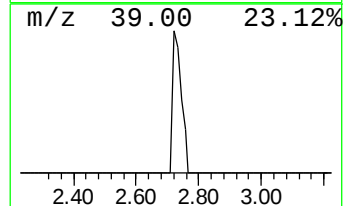
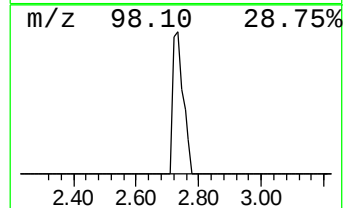
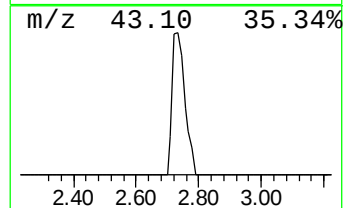
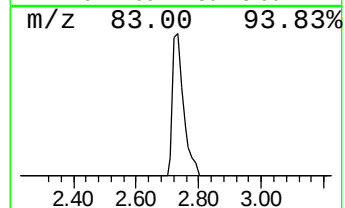
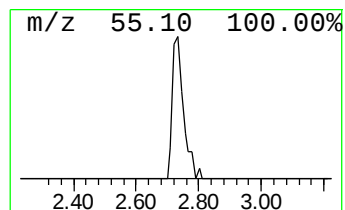
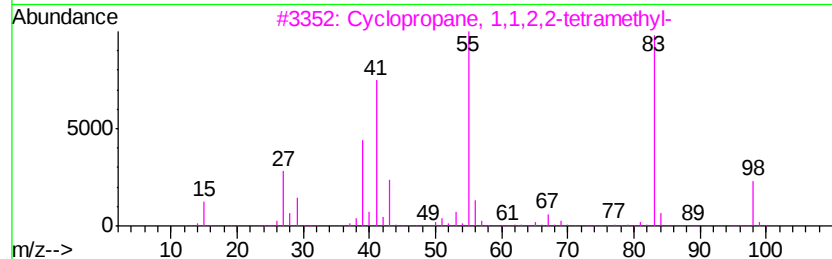
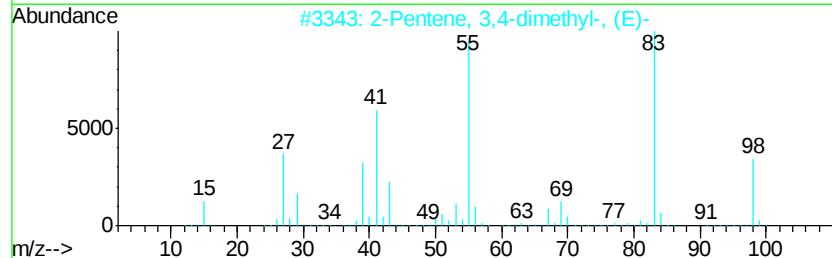
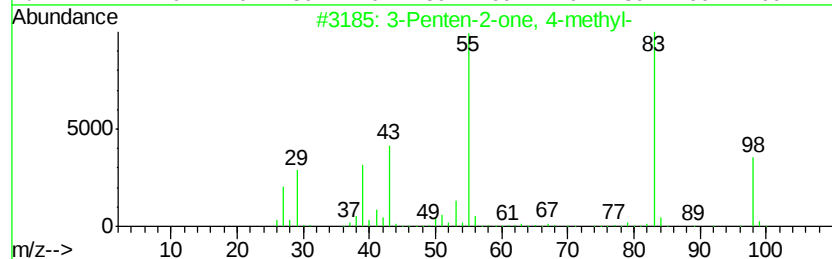
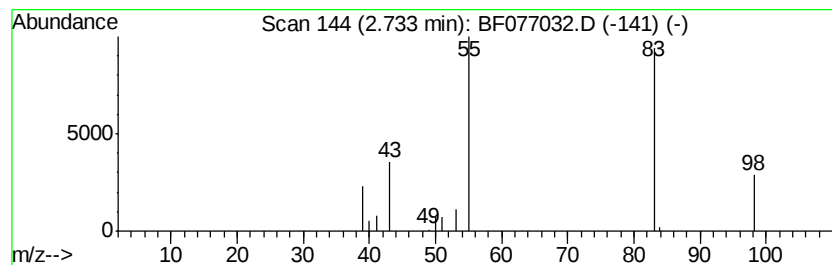
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 6

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 2.73 | 5.41 ng | 37339 | 1,4-Dichlorobenzene-d4 | 7.85 |

| Hit# | of | 5 | Tentative ID                       | MW | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|----|---------|-------------|------|
| 1    |    |   | 3-Penten-2-one, 4-methyl-          | 98 | C6H10O  | 000141-79-7 | 90   |
| 2    |    |   | 2-Pentene, 3,4-dimethyl-, (E)-     | 98 | C7H14   | 004914-92-5 | 83   |
| 3    |    |   | Cyclopropane, 1,1,2,2-tetramethyl- | 98 | C7H14   | 004127-47-3 | 78   |
| 4    |    |   | 2-Pentene, 4,4-dimethyl-, (Z)-     | 98 | C7H14   | 000762-63-0 | 78   |
| 5    |    |   | 2-Pentene, 3,4-dimethyl-, (Z)-     | 98 | C7H14   | 004914-91-4 | 78   |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

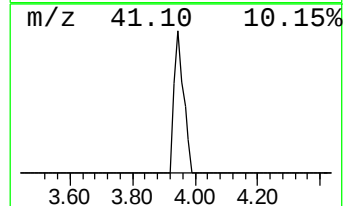
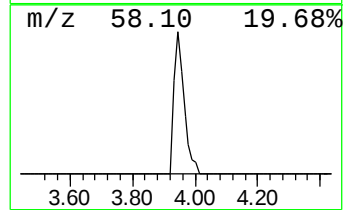
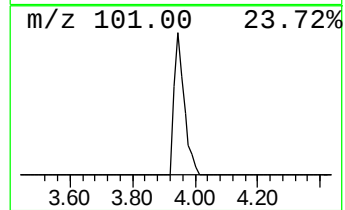
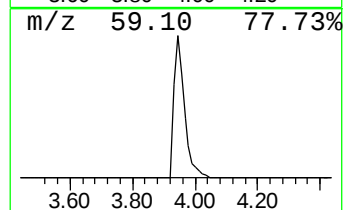
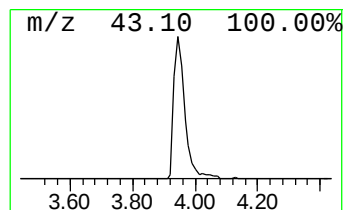
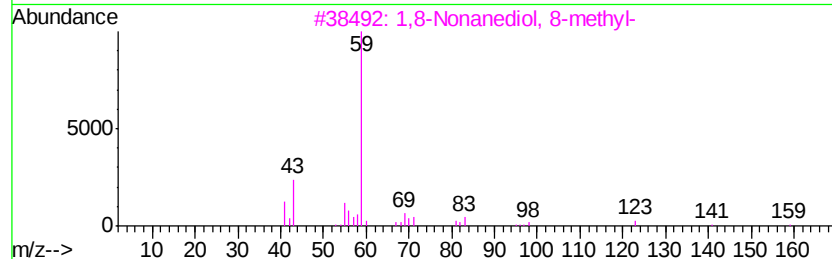
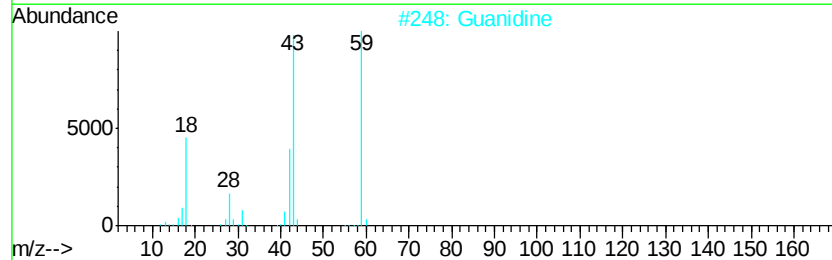
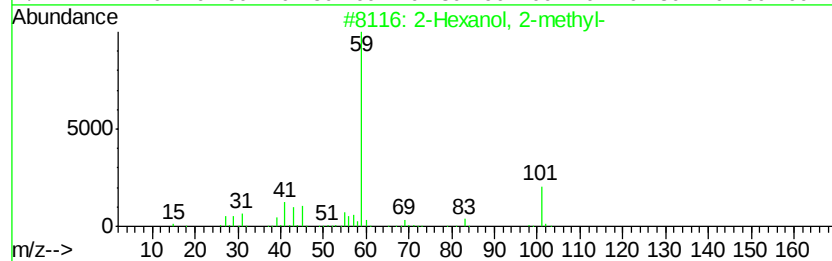
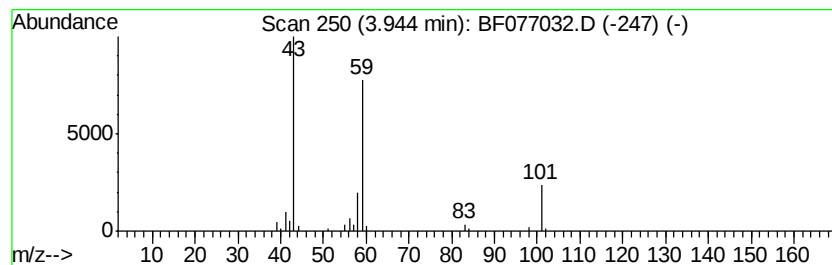
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 2-Hexanol, 2-methyl- Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 3.94 | 21.58 ng | 148862 | 1,4-Dichlorobenzene-d4 | 7.85 |

| Hit# of | 5 | Tentative ID                    | MW  | MolForm  | CAS#        | Qual |
|---------|---|---------------------------------|-----|----------|-------------|------|
| 1       |   | 2-Hexanol, 2-methyl-            | 116 | C7H16O   | 000625-23-0 | 53   |
| 2       |   | Guanidine                       | 59  | CH5N3    | 000113-00-8 | 50   |
| 3       |   | 1,8-Nonanediol, 8-methyl-       | 174 | C10H22O2 | 054725-73-4 | 38   |
| 4       |   | Butyl aldoxime, 3-methyl-, syn- | 101 | C5H11NO  | 005780-40-5 | 37   |
| 5       |   | 3-Hexanol, 4-methyl-            | 116 | C7H16O   | 000615-29-2 | 33   |



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Instrument :  
BNA\_F  
ClientSampleId :  
FK-GF-02-004-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

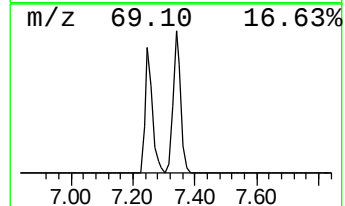
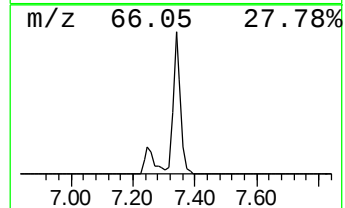
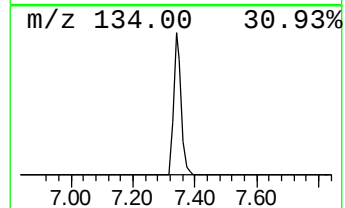
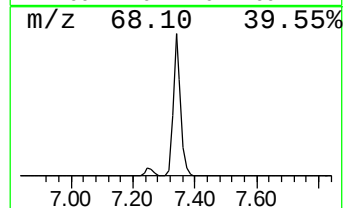
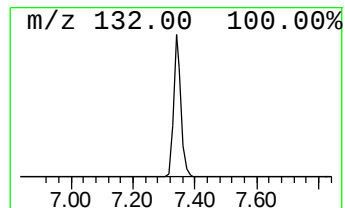
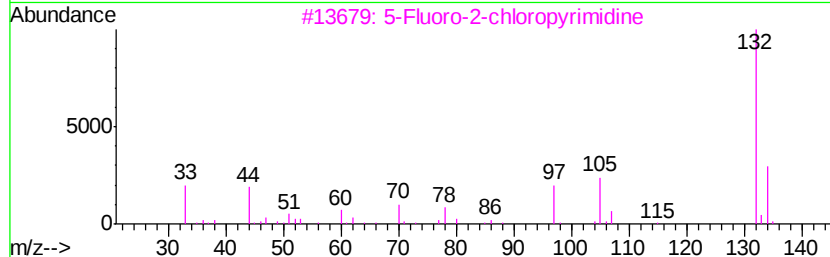
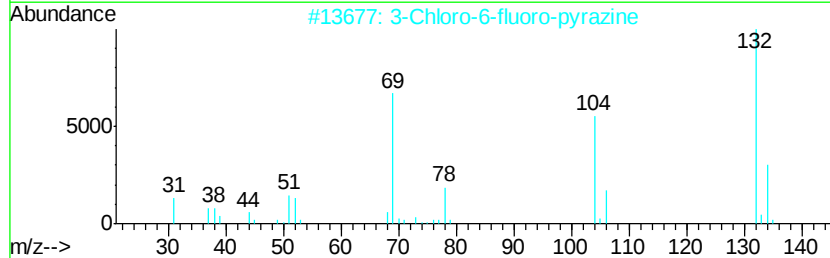
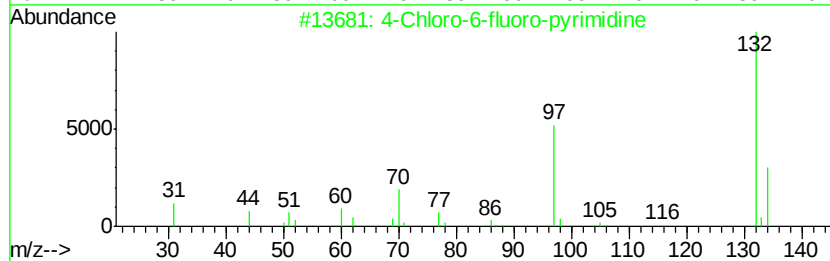
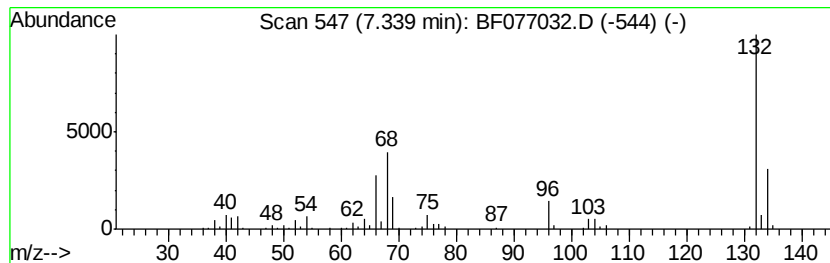
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown7.34 Concentration Rank 1

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.34 | 61.61 ng | 424939 | 1,4-Dichlorobenzene-d4 | 7.85 |

| Hit# | of | Tentative ID                 | MW  | MolForm   | CAS#         | Qual |
|------|----|------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4-Chloro-6-fluoro-pyrimidine | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 2    |    | 3-Chloro-6-fluoro-pyrazine   | 132 | C4H2ClFN2 | 1000146-10-7 | 16   |
| 3    |    | 5-Fluoro-2-chloropyrimidine  | 132 | C4H2ClFN2 | 062802-42-0  | 16   |
| 4    |    | 5-Aminoindole                | 132 | C8H8N2    | 005192-03-0  | 12   |
| 5    |    | 4-Chloro-2-fluoro-pyrimidine | 132 | C4H2ClFN2 | 051422-00-5  | 9    |



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ALS Vial : 21 Sample Multiplier: 1

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ClientSampleId :  
FK-GF-02-004-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

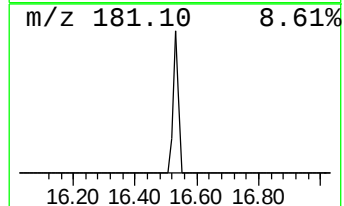
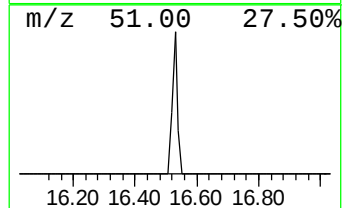
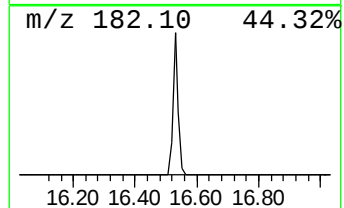
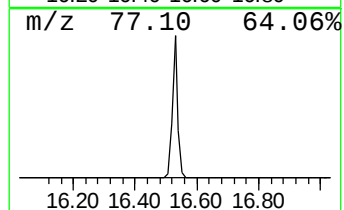
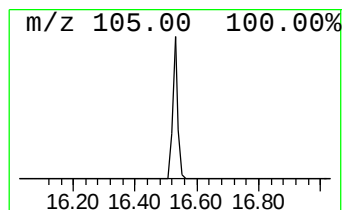
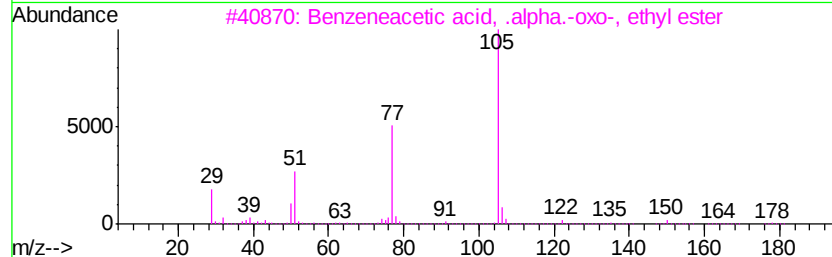
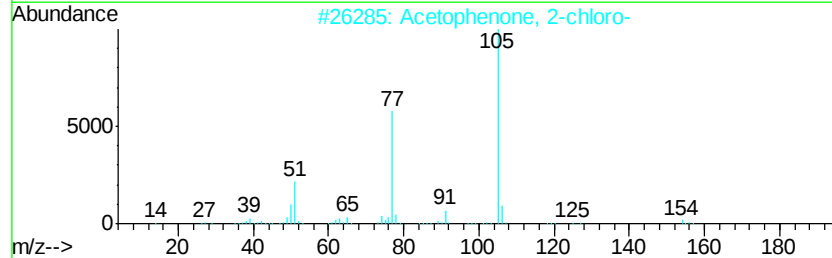
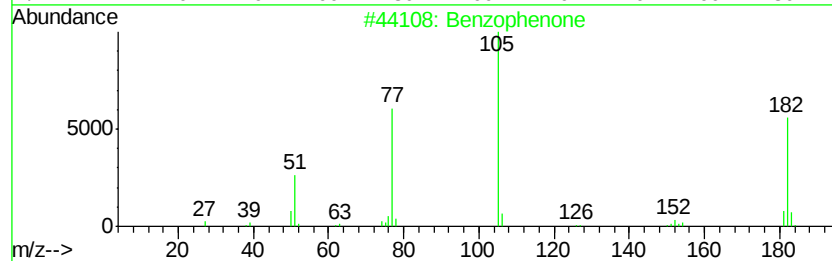
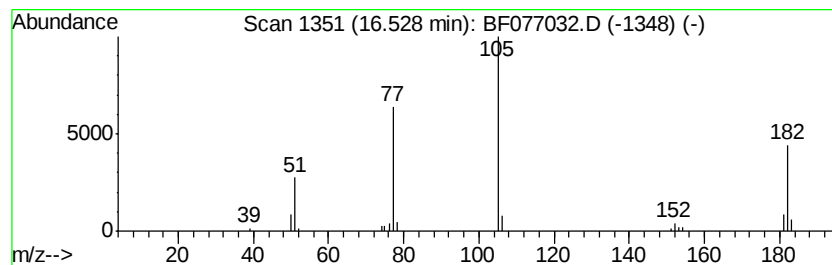
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzophenone Concentration Rank 5

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.53 | 7.47 ng | 80588 | Phenanthrene-d10 | 17.69 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|---------|---|-------------------------------------|-----|----------|--------------|------|
| 1       |   | Benzophenone                        | 182 | C13H10O  | 000119-61-9  | 96   |
| 2       |   | Acetophenone, 2-chloro-             | 154 | C8H7ClO  | 000532-27-4  | 49   |
| 3       |   | Benzeneacetic acid, .alpha.-oxo-... | 178 | C10H10O3 | 001603-79-8  | 47   |
| 4       |   | .alpha.,.alpha.-Dichloroacetophe... | 188 | C8H6Cl2O | 002648-61-5  | 47   |
| 5       |   | N-(1-Cyanovinyl)benzamide           | 172 | C10H8N2O | 1000186-19-1 | 47   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012215\  
Data File : BF077032.D  
Acq On : 23 Jan 2015 5:08  
Operator : TP/IZ  
Sample : G1148-03  
Misc : S  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
FK-GF-02-004-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

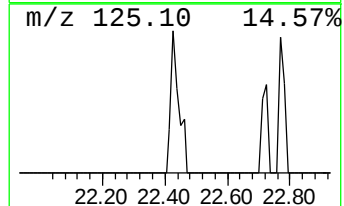
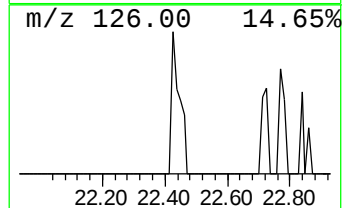
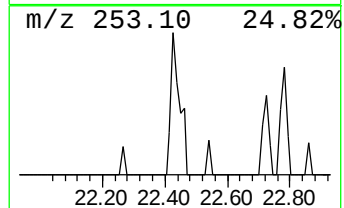
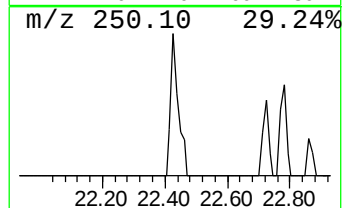
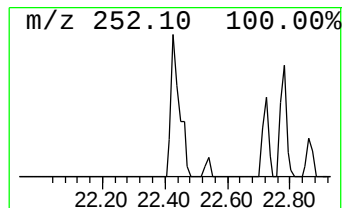
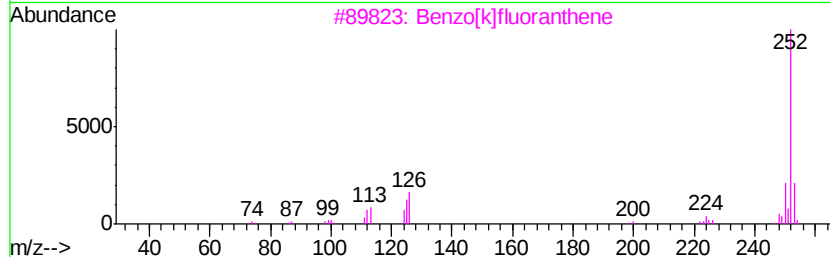
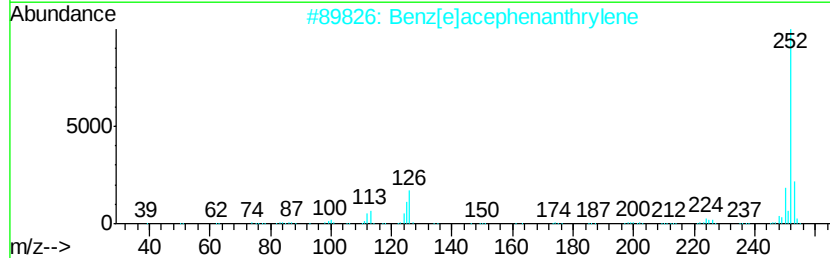
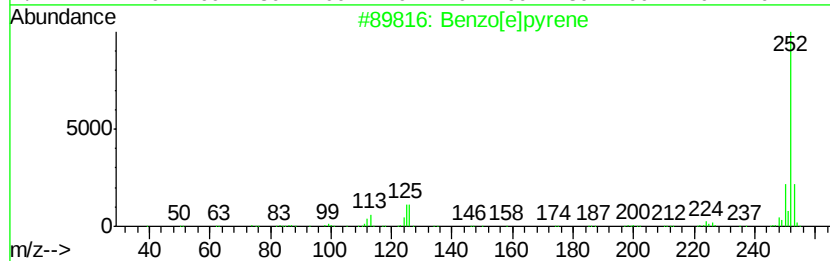
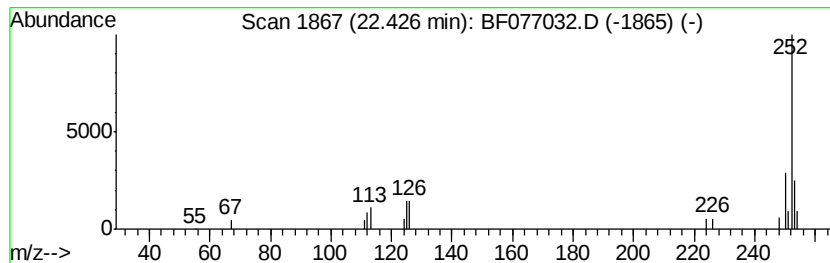
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Benzo[e]pyrene Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 22.43 | 2.69 ng | 28581 | Perylene-d12     | 22.84 |

| Hit# of 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|-----------|---------------------------|-----|---------|-------------|------|
| 1         | Benzo[e]pyrene            | 252 | C20H12  | 000192-97-2 | 87   |
| 2         | Benzo[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 87   |
| 3         | Benzo[k]fluoranthene      | 252 | C20H12  | 000207-08-9 | 81   |
| 4         | Benzo[a]pyrene            | 252 | C20H12  | 000050-32-8 | 81   |
| 5         | Perylene                  | 252 | C20H12  | 000198-55-0 | 72   |



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Data File : BF077032.D  
Acq On : 23 Jan 2015 5:08  
Operator : TP/IZ  
Sample : G1148-03  
Misc : S  
ALS Vial : 21 Sample Multiplier: 1

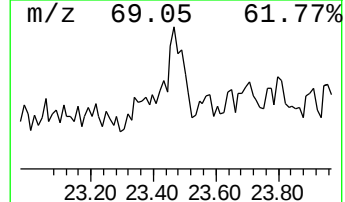
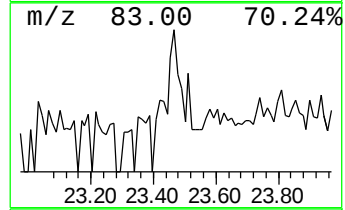
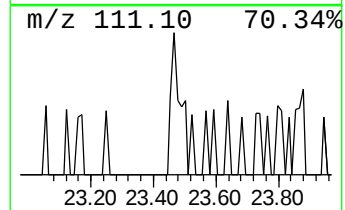
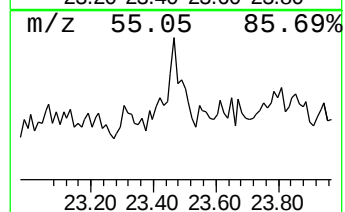
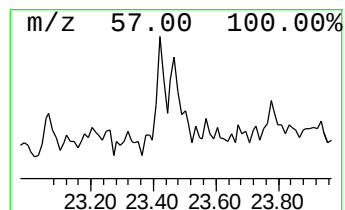
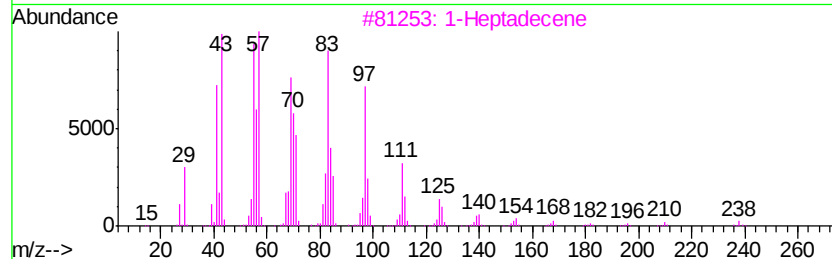
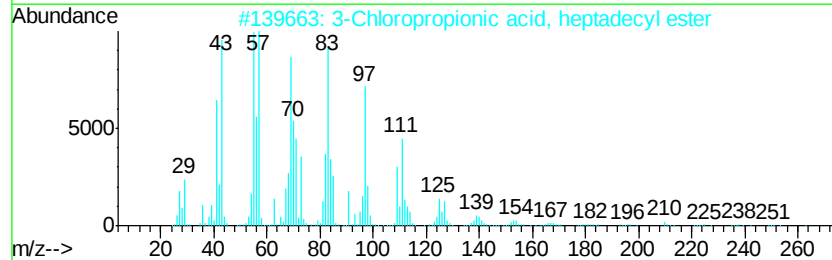
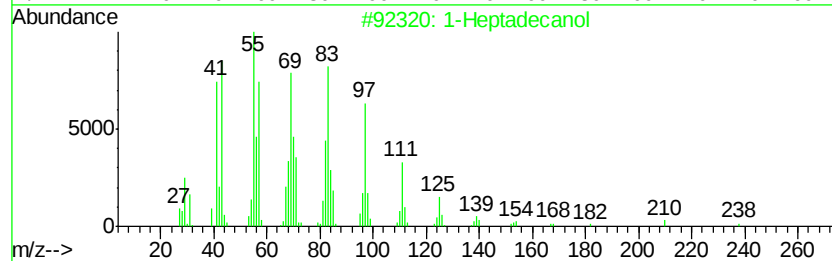
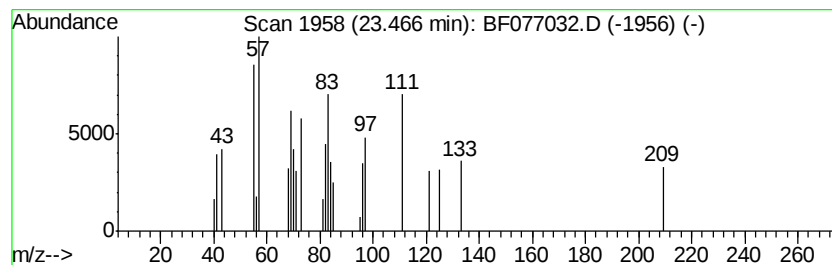
Instrument :  
BNA\_F  
ClientSampleId :  
FK-GF-02-004-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 1-Heptadecanol Concentration Rank 8

| R.T.    | EstConc | Area                                | Relative to ISTD |            | R.T.         |      |
|---------|---------|-------------------------------------|------------------|------------|--------------|------|
| 23.47   | 2.87 ng | 30519                               | Perylene-d12     |            | 22.84        |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm    | CAS#         | Qual |
| 1       |         | 1-Heptadecanol                      | 256              | C17H36O    | 001454-85-9  | 50   |
| 2       |         | 3-Chloropropionic acid, heptadec... | 346              | C20H39ClO2 | 1000283-05-1 | 50   |
| 3       |         | 1-Heptadecene                       | 238              | C17H34     | 006765-39-5  | 47   |
| 4       |         | 1-Octadecene                        | 252              | C18H36     | 000112-88-9  | 43   |
| 5       |         | 4-Heptafluorobutyryloxyhexadecane   | 438              | C20H33F7O2 | 1000282-97-2 | 43   |



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Acq On : 23 Jan 2015 5:08  
Operator : TP/IZ  
Sample : G1148-03  
Misc : S  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
FK-GF-02-004-B

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

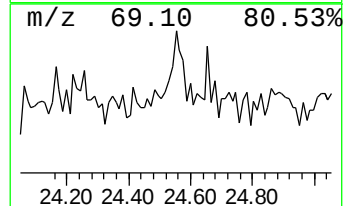
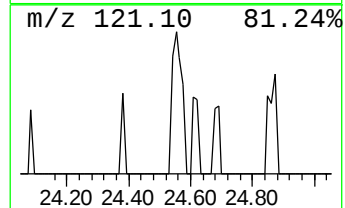
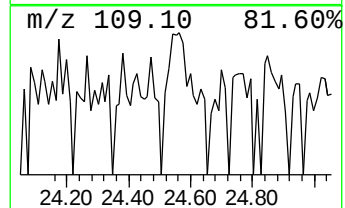
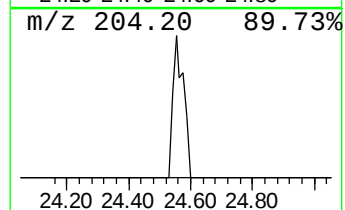
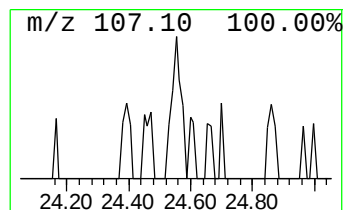
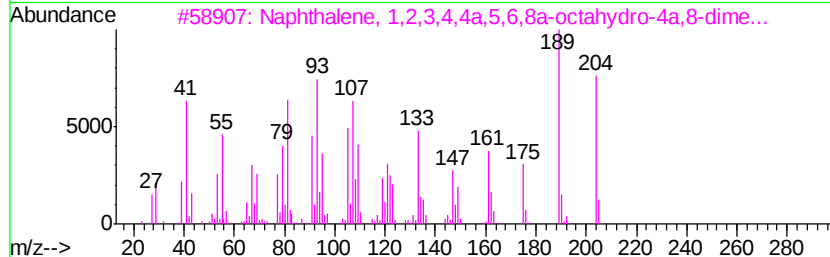
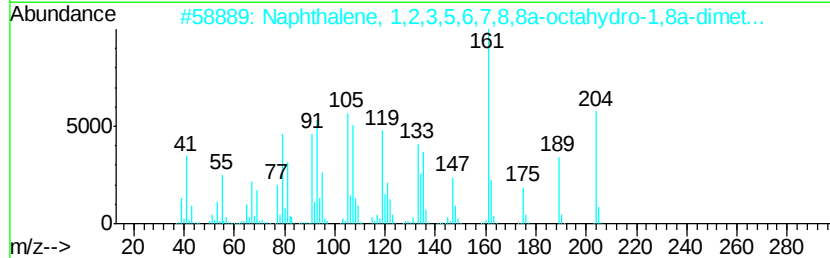
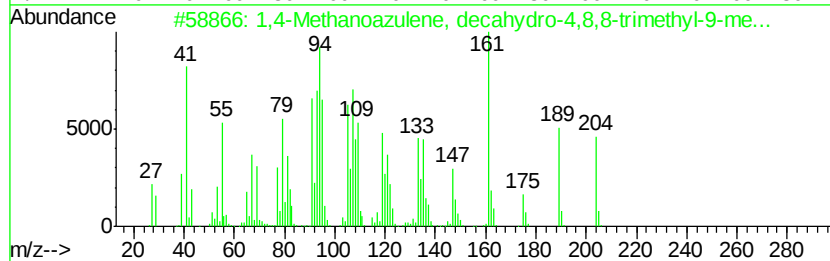
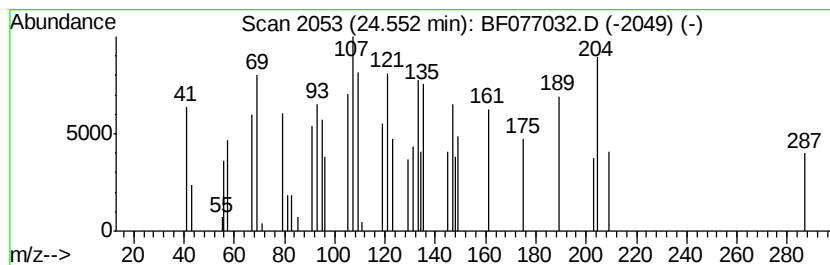
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 1,4-Methanoazulene, decahyd... Concentration Rank 7

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 24.55 | 3.41 ng | 36275 | Perylene-d12     | 22.84 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1,4-Methanoazulene, decahydro-4,... | 204 | C15H24  | 000475-20-7 | 60   |
| 2    |    |   | Naphthalene, 1,2,3,5,6,7,8,8a-oc... | 204 | C15H24  | 004630-07-3 | 60   |
| 3    |    |   | Naphthalene, 1,2,3,4,4a,5,6,8a-o... | 204 | C15H24  | 000473-13-2 | 55   |
| 4    |    |   | Naphthalene, 1,2,3,5,6,7,8,8a-oc... | 204 | C15H24  | 010219-75-7 | 53   |
| 5    |    |   | 1H-Cycloprop[e]azulene, decahydr... | 204 | C15H24  | 000489-39-4 | 50   |



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Operator : TP/IZ  
Sample : G1148-03  
Misc : S  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
FK-GF-02-004-B

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF011415.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 |
| 3-Penten-2-one, 4... | 2.73  | 5.4     | ng    | 37339    | 1                     | 7.85  | 137948 | 20.0 |
| 2-Hexanol, 2-methyl- | 3.94  | 21.6    | ng    | 148862   | 1                     | 7.85  | 137948 | 20.0 |
| unknown7.34          | 7.34  | 61.6    | ng    | 424939   | 1                     | 7.85  | 137948 | 20.0 |
| Benzophenone         | 16.53 | 7.5     | ng    | 80588    | 4                     | 17.69 | 215841 | 20.0 |
| Benzo[e]pyrene       | 22.43 | 2.7     | ng    | 28581    | 6                     | 22.84 | 212605 | 20.0 |
| 1-Heptadecanol       | 23.47 | 2.9     | ng    | 30519    | 6                     | 22.84 | 212605 | 20.0 |
| 1,4-Methanoazulen... | 24.55 | 3.4     | ng    | 36275    | 6                     | 22.84 | 212605 | 20.0 |