

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
 Data File : BF092545.D  
 Acq On : 27 Jan 2017 00:37  
 Operator : SJ/MA  
 Sample : I1451-04  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 W-35

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-BF012417.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 5.150       | 265           | 268         | 271          | rBV      | 467749         | 628088        | 12.80%          | 1.358%        |
| 2         | 5.242       | 274           | 276         | 278          | rVB      | 95144          | 91260         | 1.86%           | 0.197%        |
| 3         | 5.562       | 301           | 304         | 306          | rBV      | 2532276        | 2456411       | 50.07%          | 5.312%        |
| 4         | 6.590       | 391           | 394         | 396          | rBV      | 1141557        | 1592970       | 32.47%          | 3.445%        |
| 5         | 6.727       | 404           | 406         | 409          | rBV      | 2927459        | 3894040       | 79.37%          | 8.422%        |
| 6         | 6.967       | 424           | 427         | 429          | rBV      | 1172361        | 1021507       | 20.82%          | 2.209%        |
| 7         | 7.127       | 438           | 441         | 443          | rBV      | 3622085        | 3523439       | 71.82%          | 7.620%        |
| 8         | 7.539       | 474           | 477         | 479          | rBV      | 2614097        | 3380684       | 68.91%          | 7.311%        |
| 9         | 8.259       | 537           | 540         | 543          | rBV      | 1473878        | 1352158       | 27.56%          | 2.924%        |
| 10        | 8.499       | 558           | 561         | 563          | rBV      | 71459          | 74311         | 1.51%           | 0.161%        |
| 11        | 8.979       | 599           | 603         | 606          | rBV      | 59565          | 80674         | 1.64%           | 0.174%        |
| 12        | 9.345       | 632           | 635         | 638          | rBV      | 4622871        | 4905900       | 100.00%         | 10.610%       |
| 13        | 9.733       | 666           | 669         | 673          | rVB      | 74040          | 108292        | 2.21%           | 0.234%        |
| 14        | 9.836       | 676           | 678         | 680          | rVB2     | 111739         | 114221        | 2.33%           | 0.247%        |
| 15        | 10.031      | 692           | 695         | 697          | rBV      | 1126268        | 1525531       | 31.10%          | 3.299%        |
| 16        | 10.236      | 709           | 713         | 715          | rBV3     | 31843          | 86688         | 1.77%           | 0.187%        |
| 17        | 10.453      | 730           | 732         | 733          | rBV2     | 59218          | 49576         | 1.01%           | 0.107%        |
| 18        | 10.533      | 737           | 739         | 741          | rBV2     | 56594          | 82692         | 1.69%           | 0.179%        |
| 19        | 10.671      | 749           | 751         | 753          | rBV2     | 74605          | 108997        | 2.22%           | 0.236%        |
| 20        | 10.819      | 759           | 764         | 766          | rBV      | 2720758        | 3313834       | 67.55%          | 7.167%        |
| 21        | 10.945      | 772           | 775         | 777          | rVB      | 361664         | 361198        | 7.36%           | 0.781%        |
| 22        | 10.991      | 777           | 779         | 781          | rBV      | 396805         | 562163        | 11.46%          | 1.216%        |
| 23        | 11.025      | 781           | 782         | 784          | rVV2     | 598647         | 779499        | 15.89%          | 1.686%        |
| 24        | 11.059      | 784           | 785         | 789          | rVV      | 578946         | 902228        | 18.39%          | 1.951%        |
| 25        | 11.128      | 789           | 791         | 794          | rVV2     | 284125         | 614707        | 12.53%          | 1.329%        |
| 26        | 11.174      | 794           | 795         | 797          | rVV2     | 612669         | 629413        | 12.83%          | 1.361%        |
| 27        | 11.208      | 797           | 798         | 801          | rVV2     | 440721         | 749247        | 15.27%          | 1.620%        |
| 28        | 11.254      | 801           | 802         | 804          | rVB      | 425211         | 321720        | 6.56%           | 0.696%        |
| 29        | 11.379      | 810           | 813         | 817          | rBV3     | 76938          | 147230        | 3.00%           | 0.318%        |
| 30        | 11.516      | 822           | 825         | 827          | rBV      | 1643019        | 1578470       | 32.17%          | 3.414%        |
| 31        | 11.722      | 841           | 843         | 845          | rBV      | 202440         | 177960        | 3.63%           | 0.385%        |
| 32        | 11.802      | 848           | 850         | 854          | rVB      | 203746         | 210909        | 4.30%           | 0.456%        |
| 33        | 12.042      | 869           | 871         | 878          | rBV      | 606812         | 740111        | 15.09%          | 1.601%        |
| 34        | 12.168      | 880           | 882         | 885          | rVV      | 220865         | 220097        | 4.49%           | 0.476%        |

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Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

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-BF012417.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |         |
|----|--------|------|------|------|------|---------|---------|--------|---------|
| 35 | 12.739 | 930  | 932  | 938  | rBV2 | 82821   | 201179  | 4.10%  | 0.435%  |
| 36 | 12.831 | 938  | 940  | 943  | rVV2 | 863562  | 1064139 | 21.69% | 2.301%  |
| 37 | 12.877 | 943  | 944  | 947  | rVV  | 231502  | 286190  | 5.83%  | 0.619%  |
| 38 | 12.934 | 947  | 949  | 952  | rVB  | 83300   | 101131  | 2.06%  | 0.219%  |
| 39 | 12.991 | 952  | 954  | 959  | rVB  | 57329   | 85760   | 1.75%  | 0.185%  |
| 40 | 13.105 | 961  | 964  | 967  | rBV  | 3854910 | 4651984 | 94.82% | 10.061% |
| 41 | 13.665 | 1011 | 1013 | 1015 | rBV  | 142142  | 147418  | 3.00%  | 0.319%  |
| 42 | 13.882 | 1031 | 1032 | 1035 | rBV2 | 91796   | 117687  | 2.40%  | 0.255%  |
| 43 | 13.997 | 1040 | 1042 | 1046 | rVB  | 98572   | 111195  | 2.27%  | 0.240%  |
| 44 | 14.157 | 1053 | 1056 | 1058 | rBV  | 1529385 | 1424952 | 29.05% | 3.082%  |
| 45 | 14.774 | 1108 | 1110 | 1113 | rVB3 | 34199   | 51658   | 1.05%  | 0.112%  |
| 46 | 15.288 | 1152 | 1155 | 1160 | rVB  | 41646   | 67321   | 1.37%  | 0.146%  |
| 47 | 15.643 | 1182 | 1186 | 1188 | rBV  | 646180  | 1072271 | 21.86% | 2.319%  |
| 48 | 15.677 | 1188 | 1189 | 1194 | rVB  | 71067   | 76740   | 1.56%  | 0.166%  |
| 49 | 16.123 | 1225 | 1228 | 1231 | rBV  | 56764   | 96506   | 1.97%  | 0.209%  |
| 50 | 16.637 | 1270 | 1273 | 1281 | rVB  | 64332   | 142368  | 2.90%  | 0.308%  |
| 51 | 17.254 | 1323 | 1327 | 1329 | rBV  | 30577   | 69020   | 1.41%  | 0.149%  |
| 52 | 17.974 | 1387 | 1390 | 1395 | rVB3 | 30666   | 84717   | 1.73%  | 0.183%  |

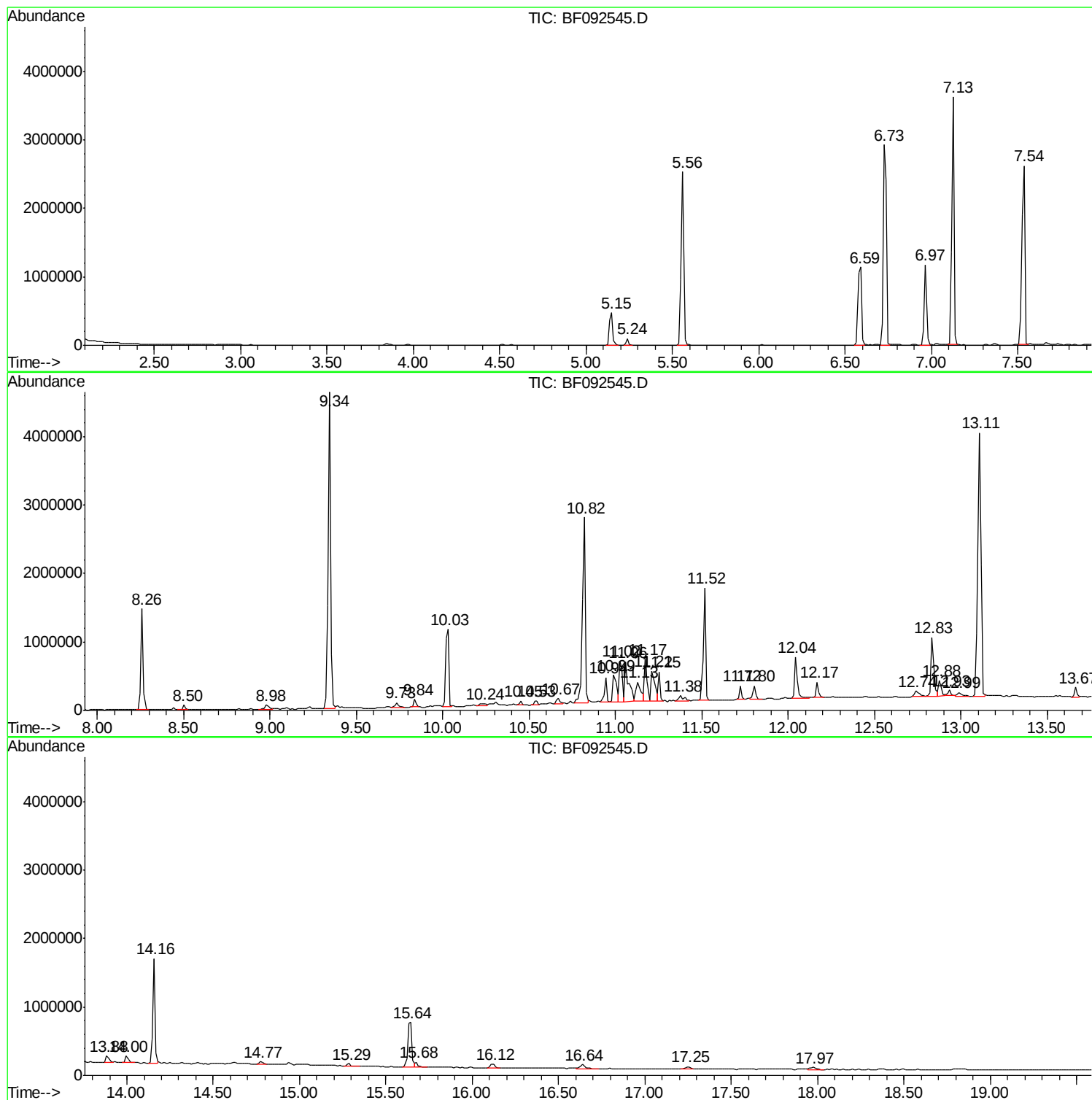
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Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
W-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012417.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\BF012617\  
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Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

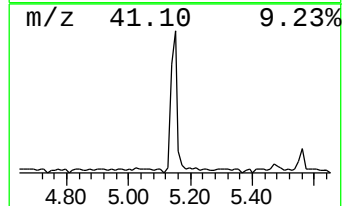
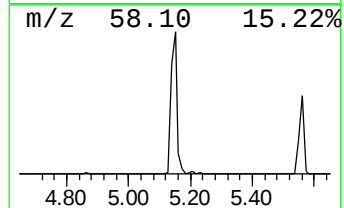
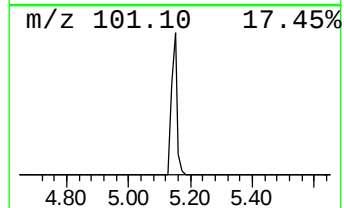
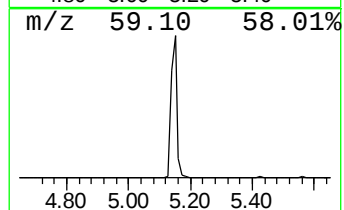
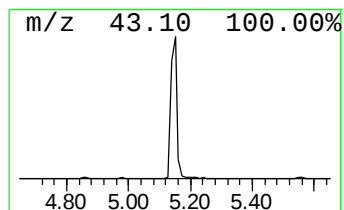
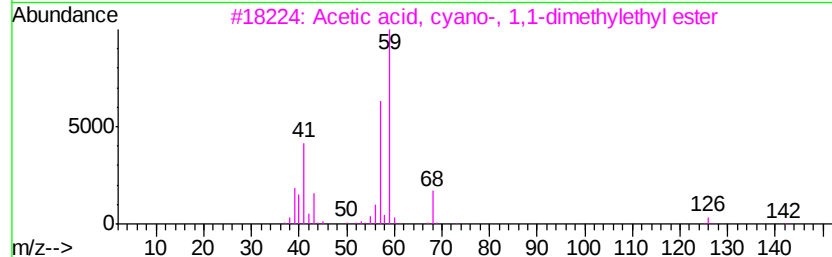
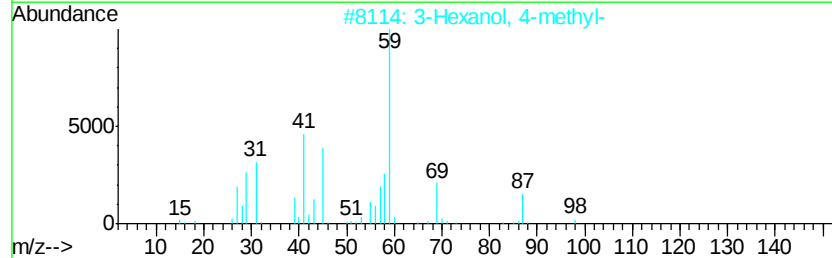
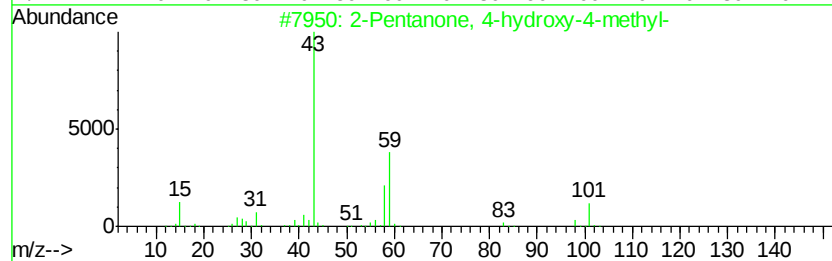
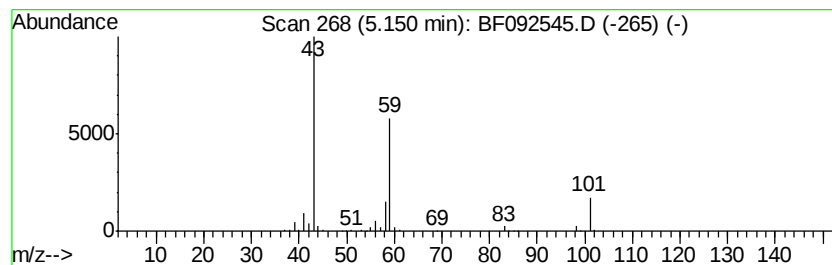
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.15 | 12.30 ng | 628088 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 45   |
| 2    |    |   | 3-Hexanol, 4-methyl-                 | 116 | C7H16O   | 000615-29-2 | 33   |
| 3    |    |   | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 25   |
| 4    |    |   | 1-Propen-2-ol, acetate               | 100 | C5H8O2   | 000108-22-5 | 10   |
| 5    |    |   | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 9    |



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Instrument :  
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
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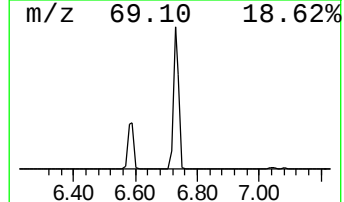
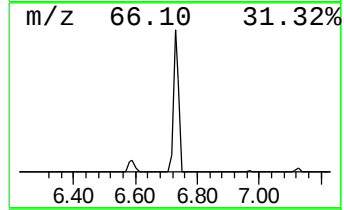
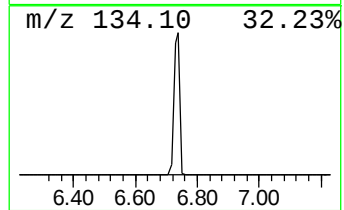
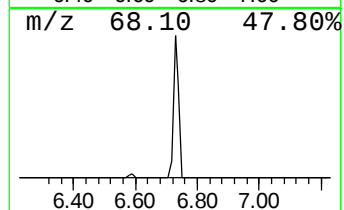
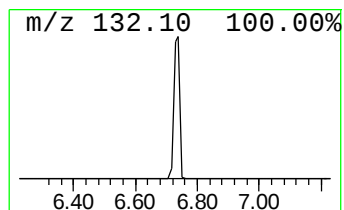
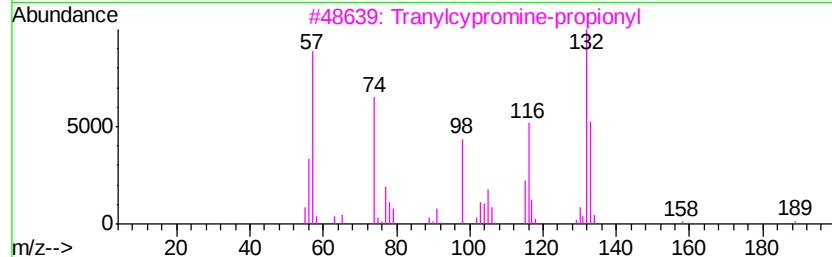
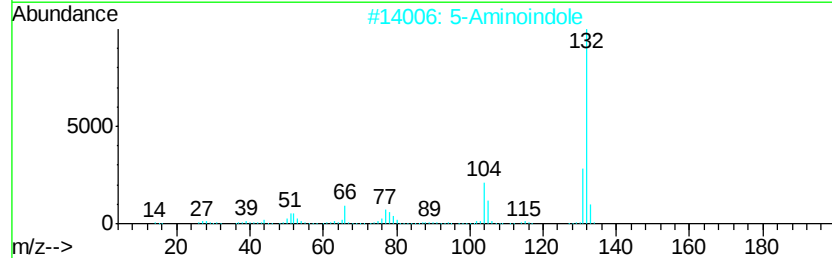
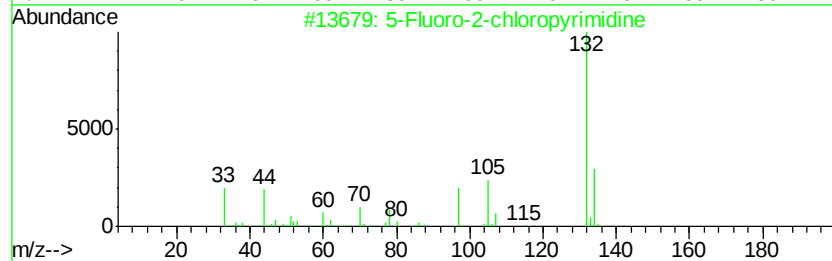
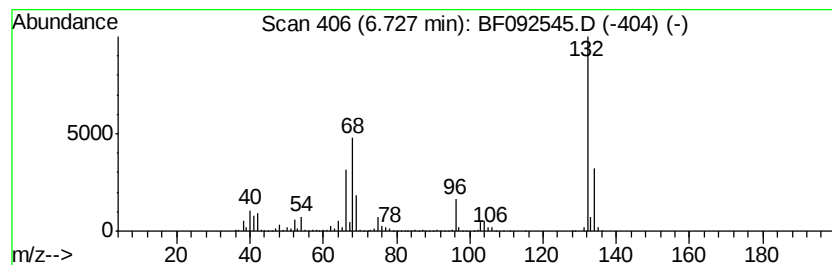
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown6.73 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.73 | 76.24 ng | 3894040 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0  | 16   |
| 2    |    | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 12   |
| 3    |    | Tranvlcyproline-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 4    |    | Benzene, 1,3,5-trifluoro-           | 132 | C6H3F3    | 000372-38-3  | 9    |
| 5    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |



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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

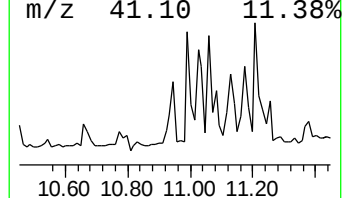
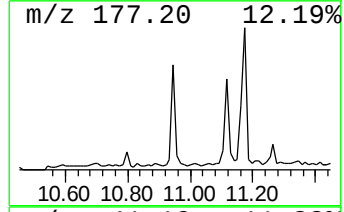
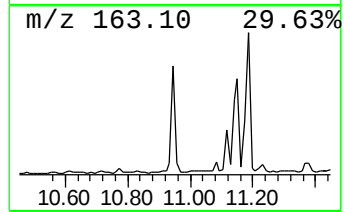
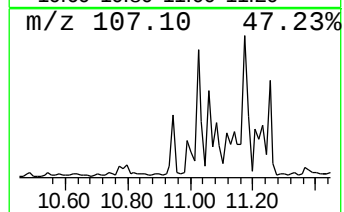
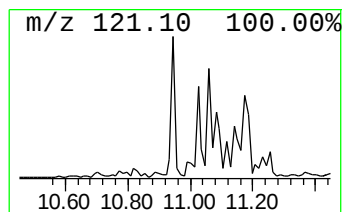
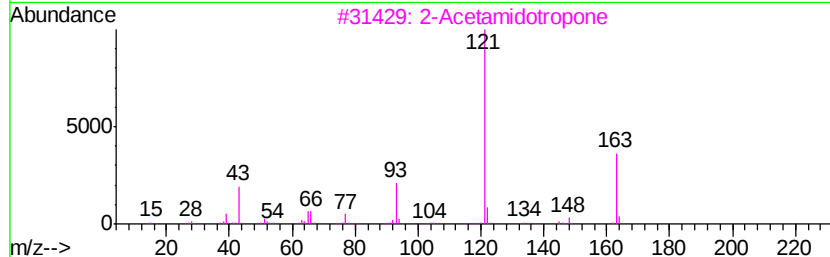
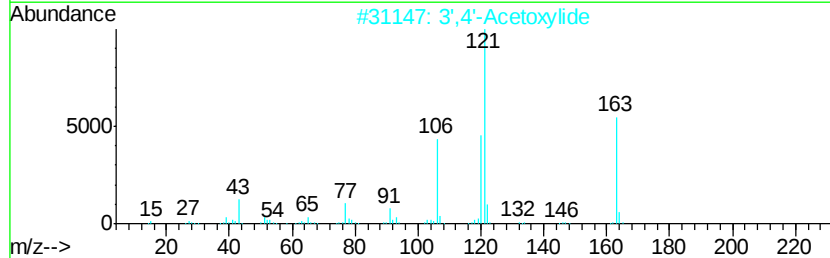
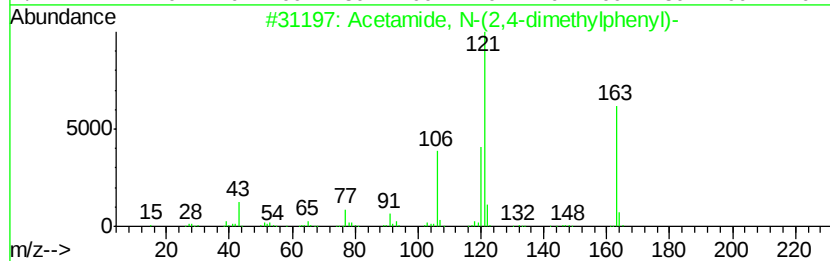
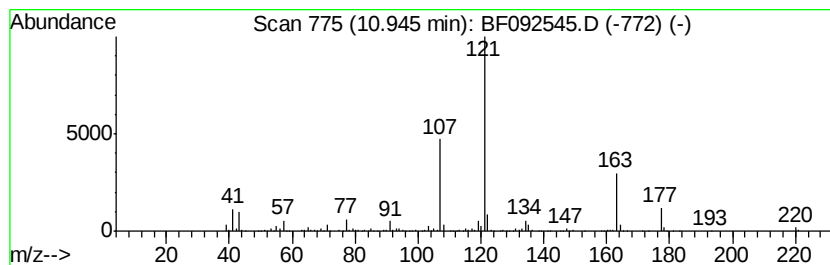
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Acetamide, N-(2,4-dimethylp... Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.95 | 4.58 ng | 361198 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Acetamide, N-(2,4-dimethylphenyl)-   | 163 | C10H13NO | 002050-43-3 | 52   |
| 2    |    | 3',4'-Acetoxylide                    | 163 | C10H13NO | 002198-54-1 | 52   |
| 3    |    | 2-Acetamidotropone                   | 163 | C9H9NO2  | 006422-12-4 | 47   |
| 4    |    | 1,2-Bis(p-acetoxylphenyl)ethanedione | 326 | C18H14O6 | 093655-79-9 | 47   |
| 5    |    | p-Sec-butylphenyl acetate            | 192 | C12H16O2 | 003245-24-7 | 43   |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

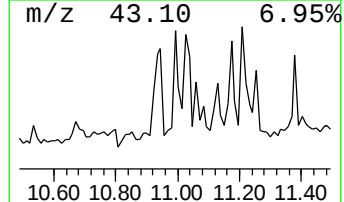
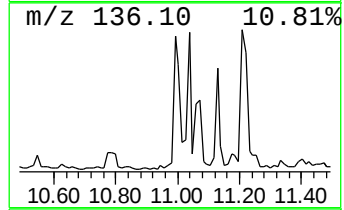
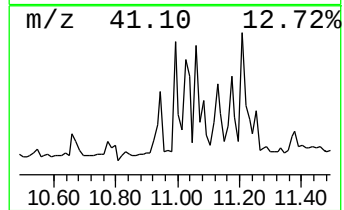
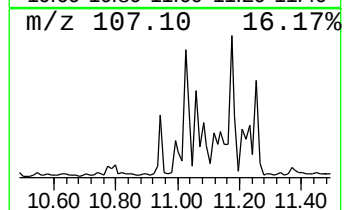
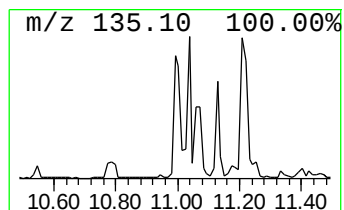
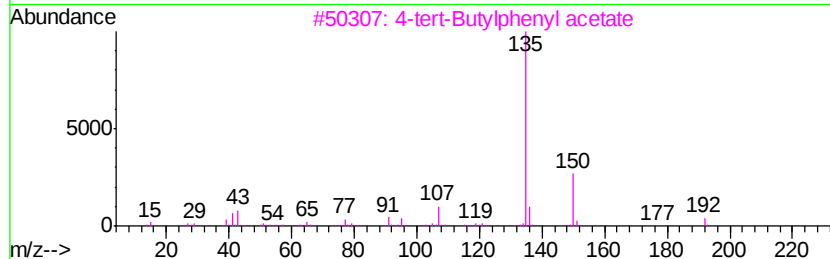
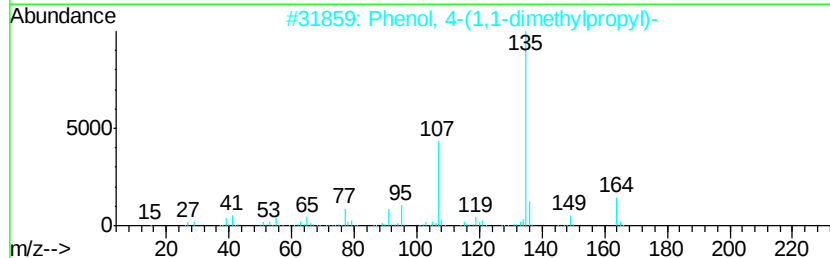
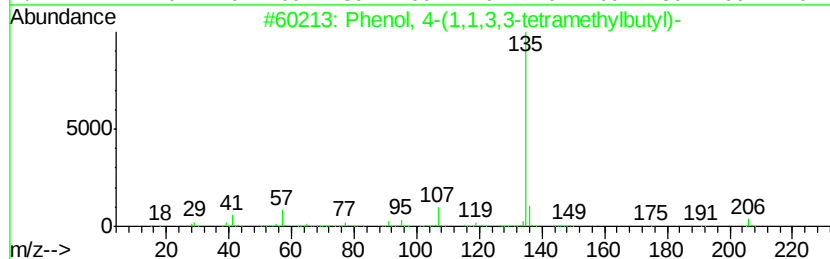
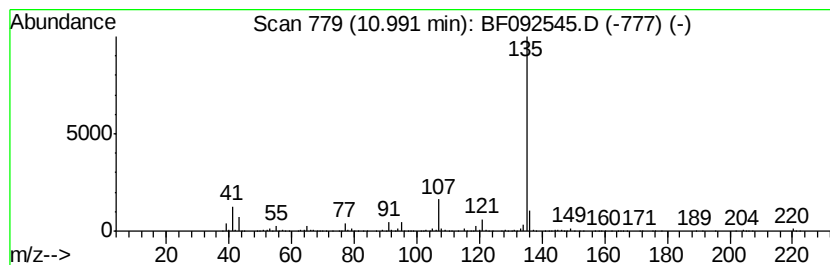
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.99 | 7.12 ng | 562163 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O     | 000140-66-9 | 78   |
| 2    |    | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O     | 000080-46-6 | 72   |
| 3    |    | 4-tert-Butylphenyl acetate          | 192 | C12H16O2    | 003056-64-2 | 72   |
| 4    |    | Pentanoic acid, 5-hydroxy-, p-t-... | 250 | C15H22O3    | 166273-37-6 | 72   |
| 5    |    | Carbamic acid, N-[1,1-bis(triflu... | 413 | C19H25F6N02 | 296242-69-8 | 56   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

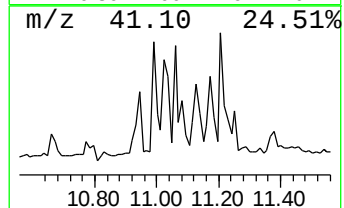
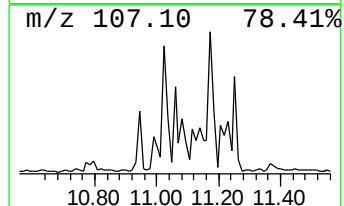
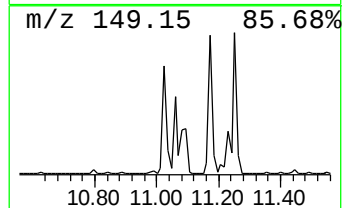
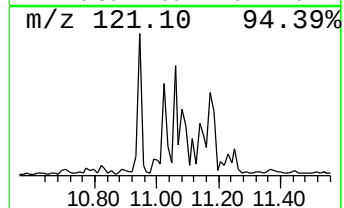
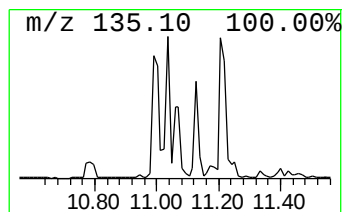
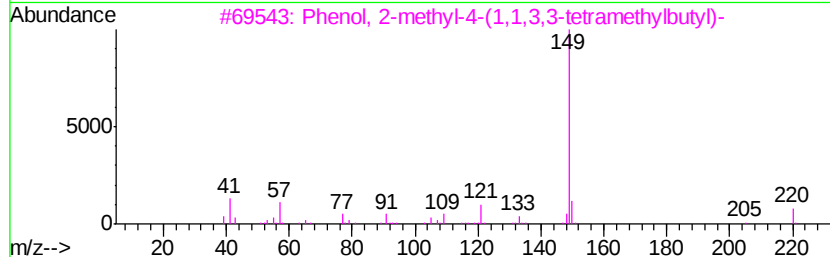
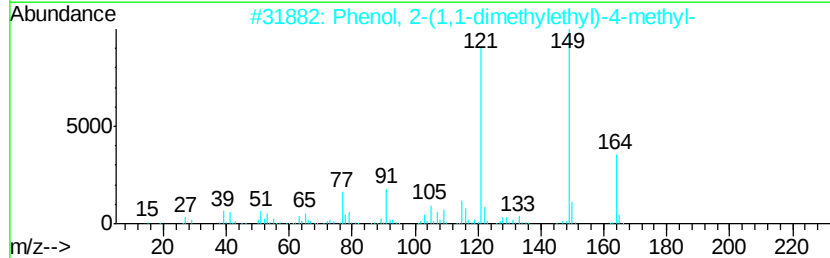
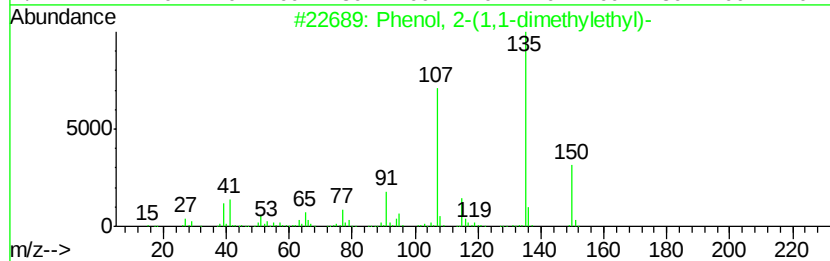
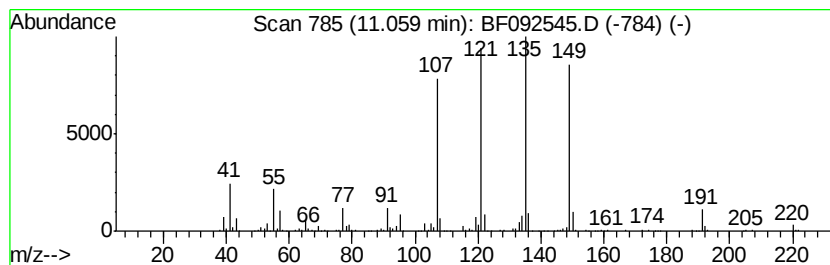
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 unknown11.06 Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.06 | 11.43 ng | 902228 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Phenol, 2-(1,1-dimethylethyl)-      | 150 | C10H14O   | 000088-18-6 | 43   |
| 2    |    | Phenol, 2-(1,1-dimethylethyl)-4-... | 164 | C11H16O   | 002409-55-4 | 35   |
| 3    |    | Phenol, 2-methyl-4-(1,1,3,3-tetr... | 220 | C15H24O   | 002219-84-3 | 18   |
| 4    |    | 4-(3,4-Methylenedioxyphenyl)-2-b... | 192 | C11H12O3  | 055418-52-5 | 14   |
| 5    |    | Phenol, 3-methyl-5-(1-methylethy... | 207 | C12H17NO2 | 002631-37-0 | 14   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

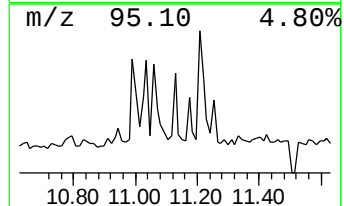
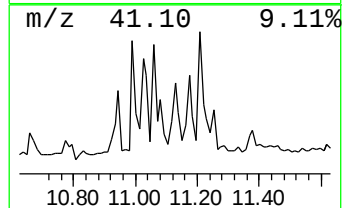
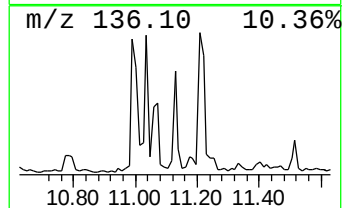
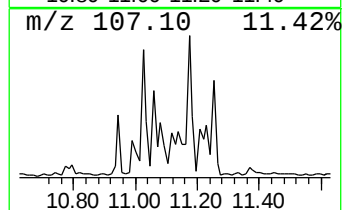
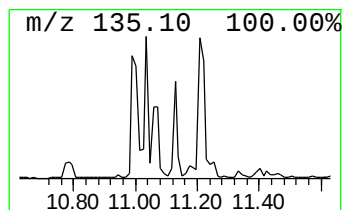
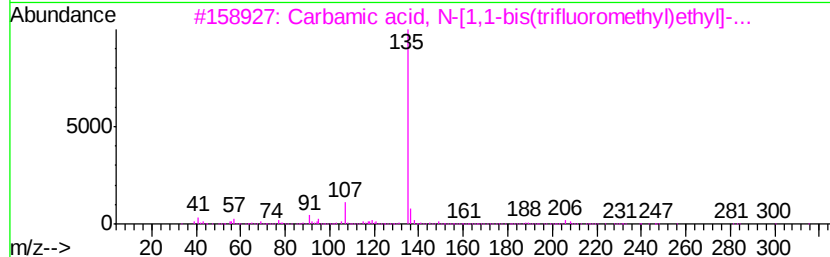
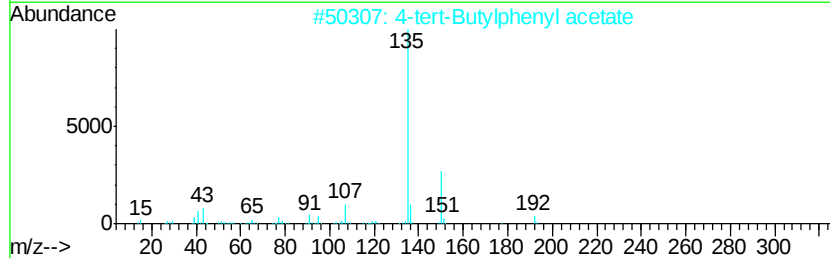
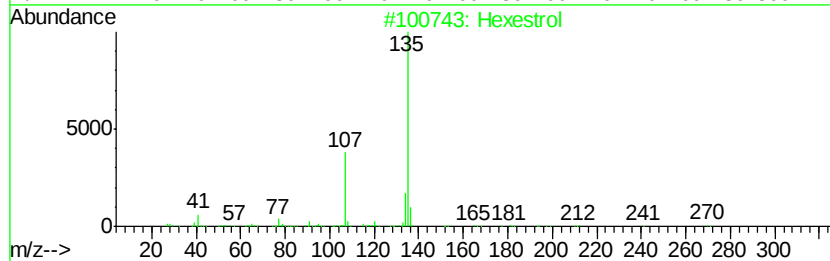
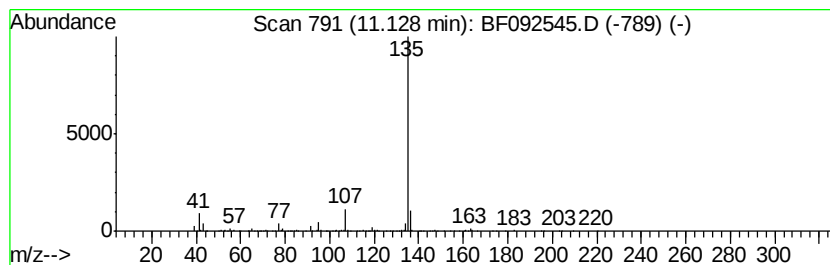
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown11.13 Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.13 | 7.79 ng | 614707 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Hexestrol                           | 270 | C18H22O2    | 005635-50-7  | 45   |
| 2    |    | 4-tert-Butylphenyl acetate          | 192 | C12H16O2    | 003056-64-2  | 43   |
| 3    |    | Carbamic acid, N-[1,1-bis(triflu... | 413 | C19H25F6N02 | 296242-69-8  | 40   |
| 4    |    | m-Methoxybenzoic acid, 2-bromo-4... | 324 | C14H10BrF03 | 1000292-63-0 | 39   |
| 5    |    | Phenol, 4,4'-(1,2-diethyl-1,2-et... | 270 | C18H22O2    | 000084-16-2  | 39   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

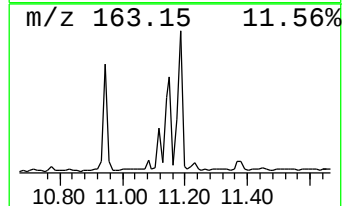
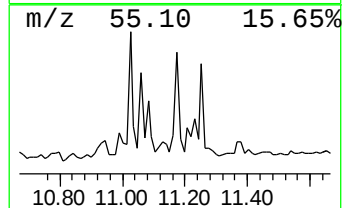
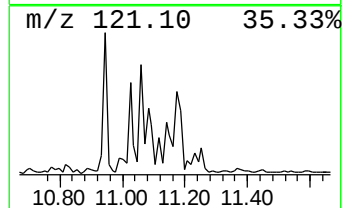
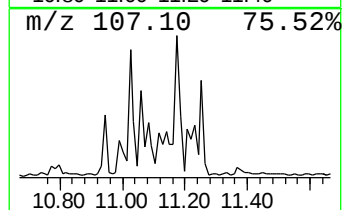
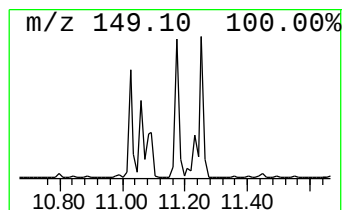
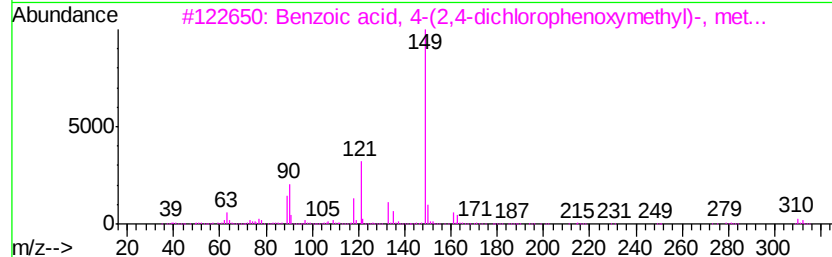
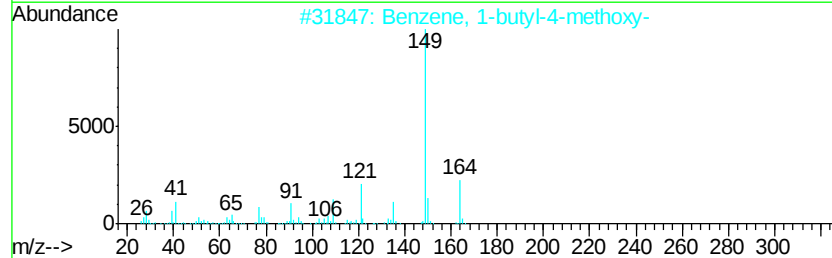
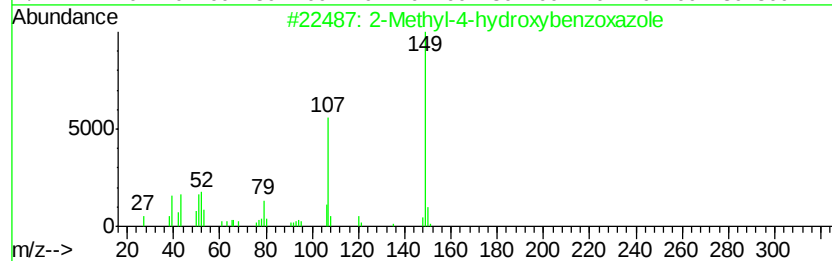
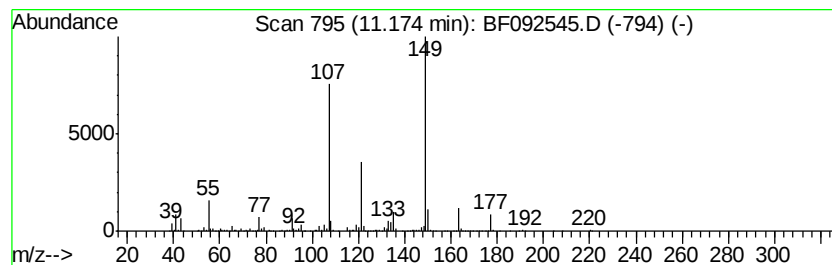
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 2-Methyl-4-hydroxybenzoxazole Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD |  | R.T.  |
|-------|---------|--------|------------------|--|-------|
| 11.17 | 7.97 ng | 629413 | Phenanthrene-d10 |  | 11.52 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 2-Methyl-4-hydroxybenzoxazole       | 149 | C8H7NO2     | 051110-60-2  | 64   |
| 2    |    |   | Benzene, 1-butyl-4-methoxy-         | 164 | C11H16O     | 018272-84-9  | 38   |
| 3    |    |   | Benzoic acid, 4-(2,4-dichlorophe... | 310 | C15H12Cl2O3 | 1000294-38-1 | 37   |
| 4    |    |   | 2H-Indol-2-one, 1,3-dihydro-5-hv... | 149 | C8H7NO2     | 003416-18-0  | 35   |
| 5    |    |   | 4-(p-Acetoxyphenyl)-2-butanone      | 206 | C12H14O3    | 003572-06-3  | 35   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

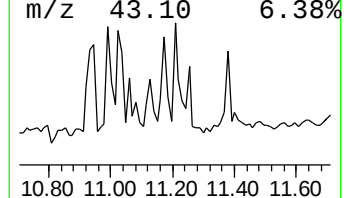
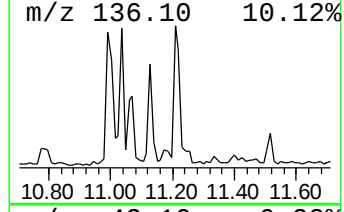
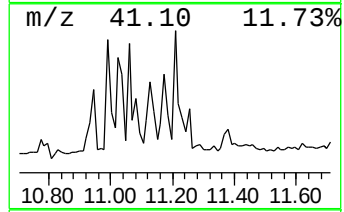
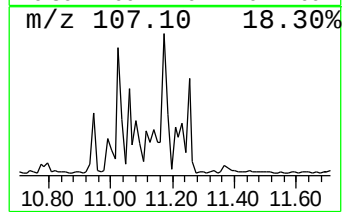
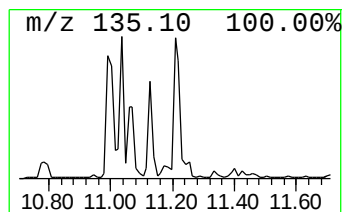
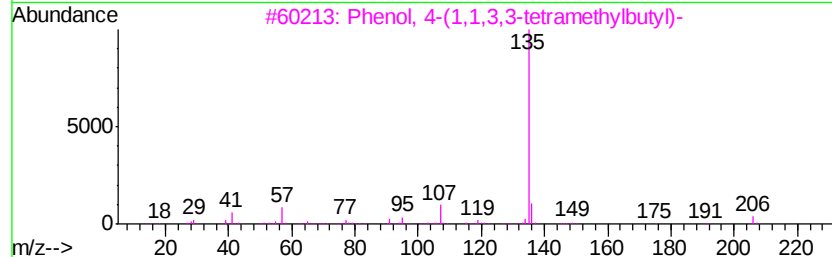
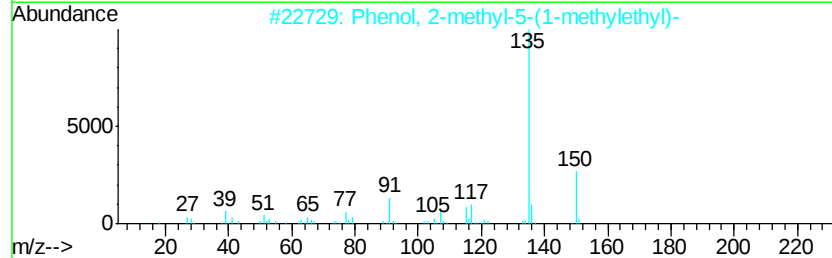
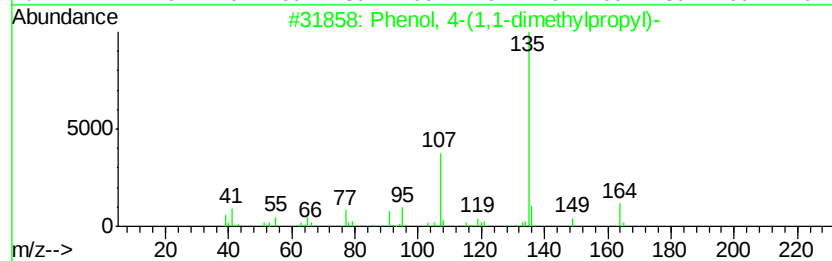
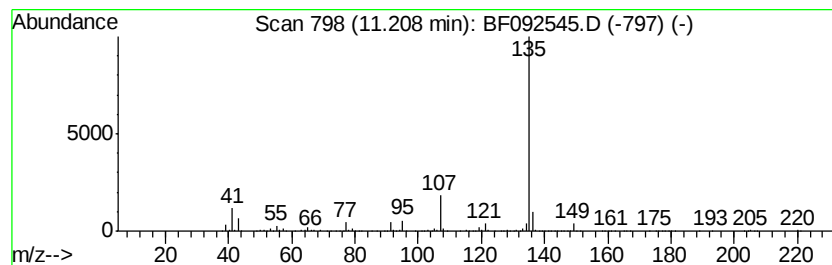
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Phenol, 4-(1,1-dimethylprop... Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.21 | 9.49 ng | 749247 | Phenanthrene-d10 | 11.52 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O  | 000080-46-6 | 72   |
| 2    |    |   | Phenol, 2-methyl-5-(1-methylethyl)- | 150 | C10H14O  | 000499-75-2 | 72   |
| 3    |    |   | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O  | 000140-66-9 | 72   |
| 4    |    |   | Hexestrol                           | 270 | C18H22O2 | 005635-50-7 | 64   |
| 5    |    |   | Pentandioic acid, (p-t-butylphen... | 264 | C15H20O4 | 212762-88-4 | 64   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

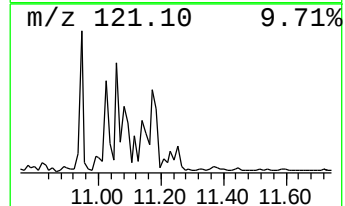
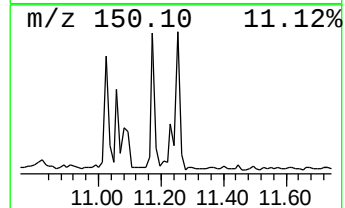
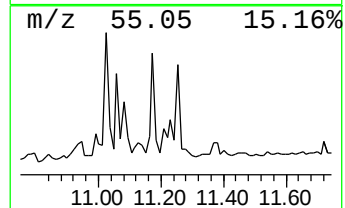
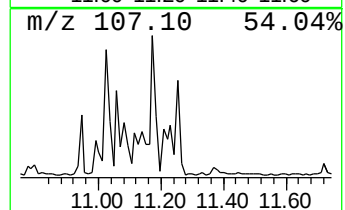
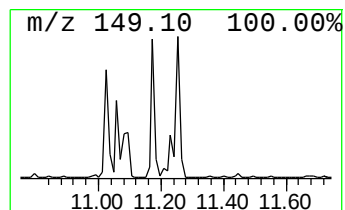
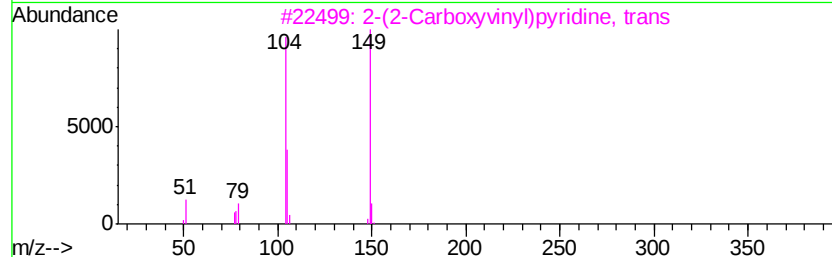
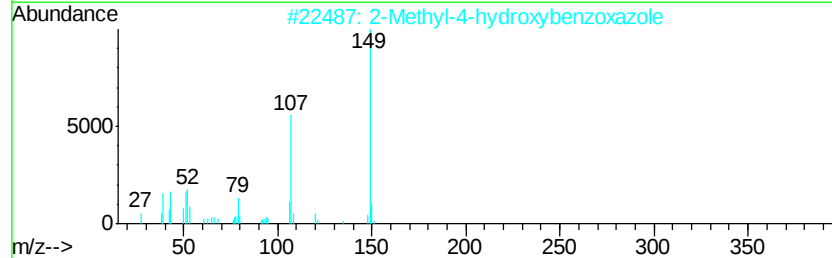
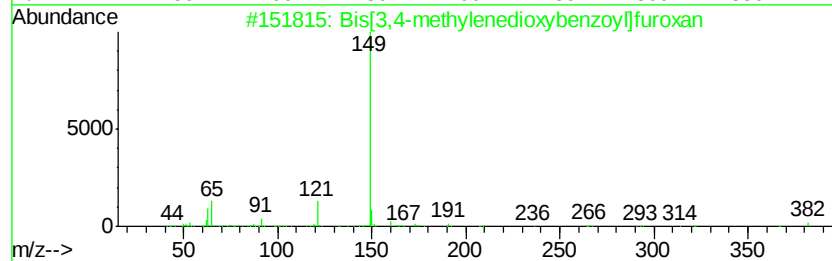
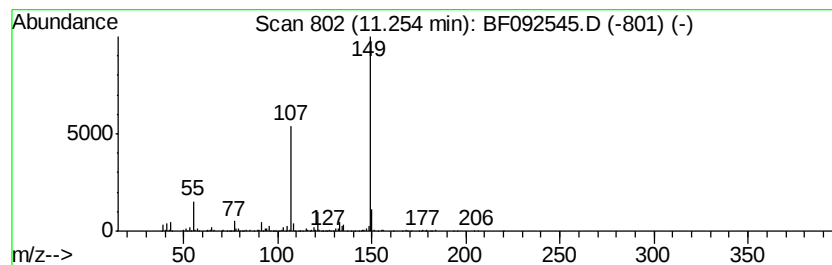
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown11.25 Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.25 | 4.08 ng | 321720 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Bis[3,4-methylenedioxybenzoyl]fu... | 382 | C18H10N2O8 | 240142-32-9  | 47   |
| 2    |    | 2-Methyl-4-hydroxybenzoxazole       | 149 | C8H7NO2    | 051110-60-2  | 39   |
| 3    |    | 2-(2-Carboxyvinyl)pyridine, trans   | 149 | C8H7NO2    | 007340-22-9  | 37   |
| 4    |    | Acetamide, 2-(4-cyclohexylphenox... | 324 | C20H24N2O2 | 1000263-95-6 | 33   |
| 5    |    | 2H-1,4-Benzoxazin-3(4H)-one         | 149 | C8H7NO2    | 005466-88-6  | 28   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

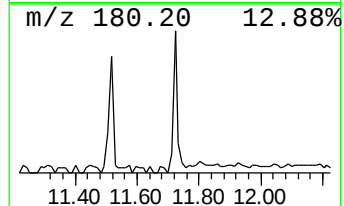
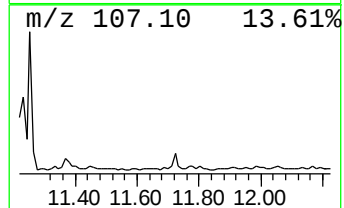
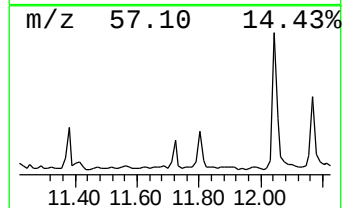
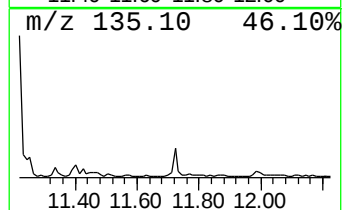
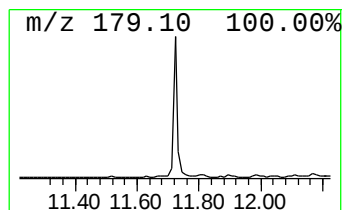
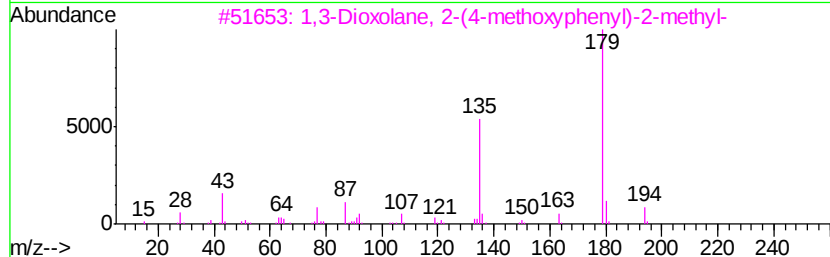
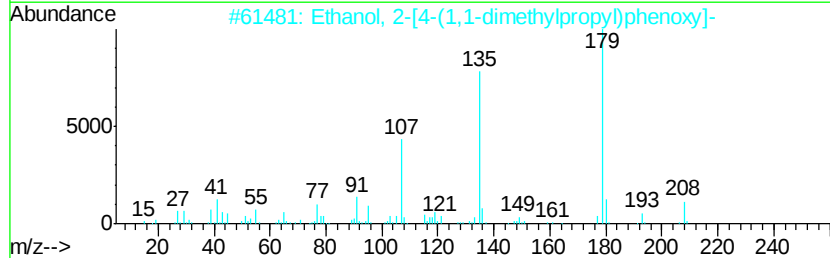
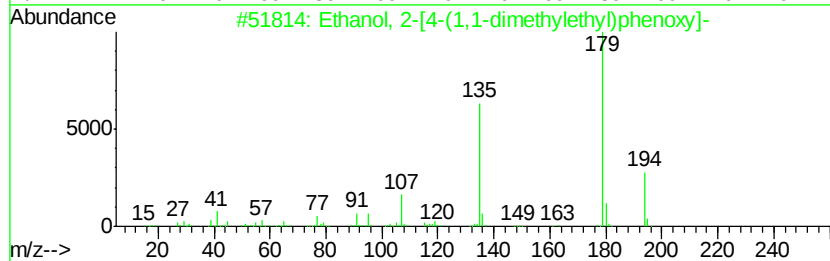
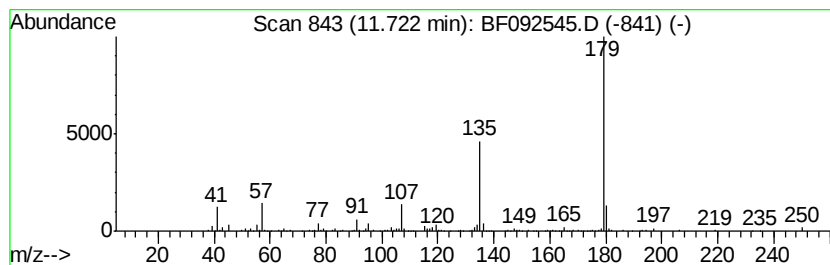
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Ethanol, 2-[4-(1,1-dimethyl... Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.72 | 2.25 ng | 177960 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                                 | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Ethanol, 2-[4-(1,1-dimethylethyl)phenoxy]-   | 194 | C12H18O2  | 000713-46-2  | 86   |
| 2    |    | Ethanol, 2-[4-(1,1-dimethylpropyl)phenoxy]-  | 208 | C13H20O2  | 006382-07-6  | 72   |
| 3    |    | 1,3-Dioxolane, 2-(4-methoxyphenyl)-2-methyl- | 194 | C11H14O3  | 036881-00-2  | 64   |
| 4    |    | 6H-Purin-6-one, 2-(dimethylamino)-           | 179 | C7H9N5O   | 001445-15-4  | 38   |
| 5    |    | 1-Cyclohexyldimethylsilyloxy-2-p-            | 262 | C16H26OSi | 1000282-21-3 | 25   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

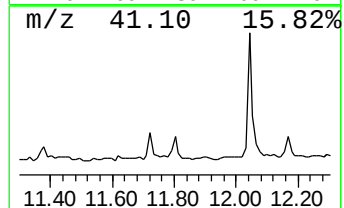
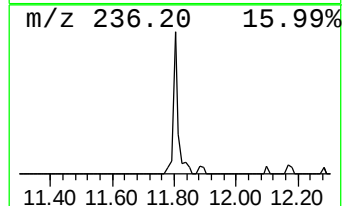
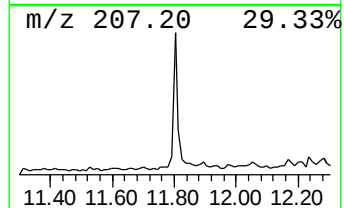
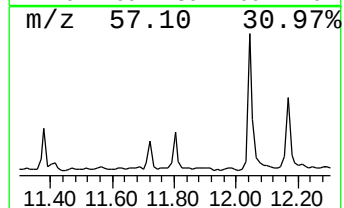
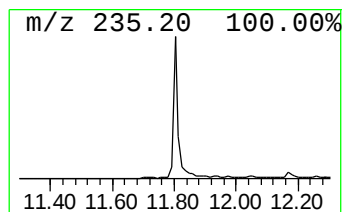
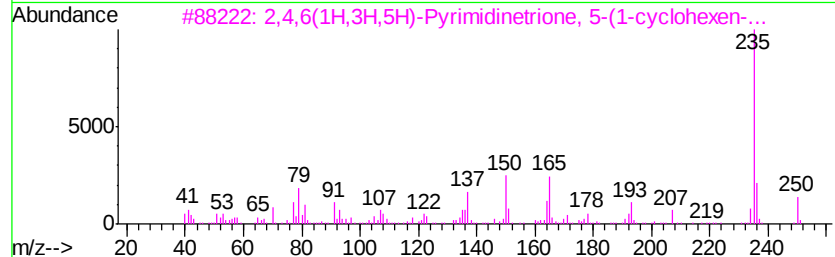
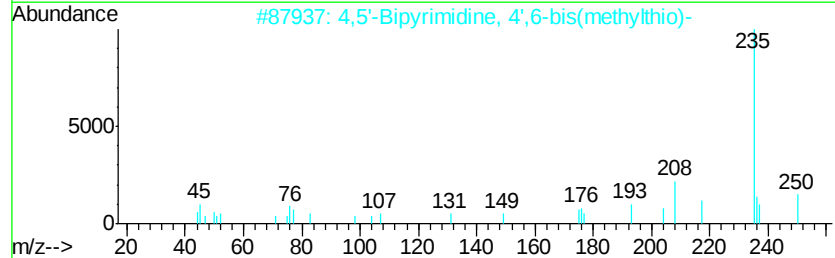
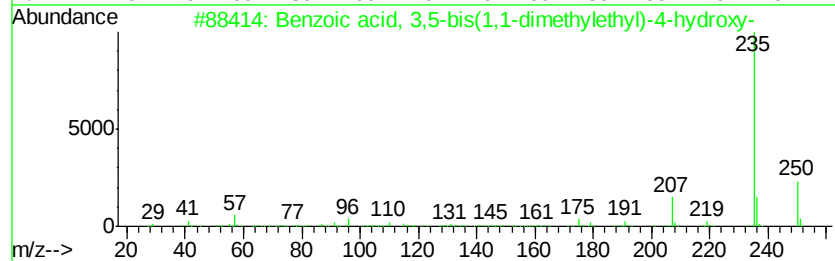
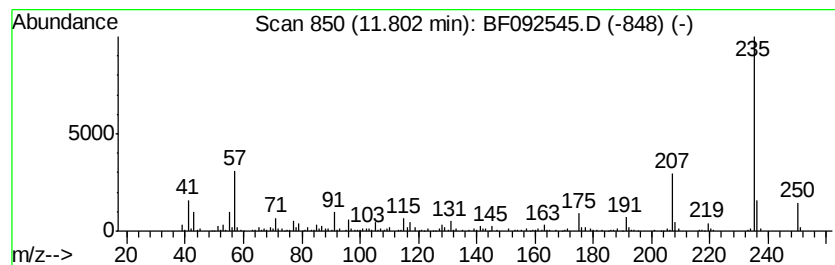
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Benzoic acid, 3,5-bis(1,1-d... Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.80 | 2.67 ng | 210909 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Benzoic acid, 3,5-bis(1,1-dimeth... | 250 | C15H22O3   | 001421-49-4  | 94   |
| 2    |    | 4,5'-Bipyriridine, 4',6-bis(meth... | 250 | C10H10N4S2 | 059549-36-9  | 59   |
| 3    |    | 2,4,6(1H,3H,5H)-Pyrimidinetrione... | 250 | C13H18N2O3 | 000726-79-4  | 53   |
| 4    |    | N-(4-Hydroxyphenyl)-2-naphthylamine | 235 | C16H13NO   | 000093-45-8  | 53   |
| 5    |    | Hexobarbital-2(or 4)-methyl deri... | 250 | C13H18N2O3 | 1000123-51-5 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

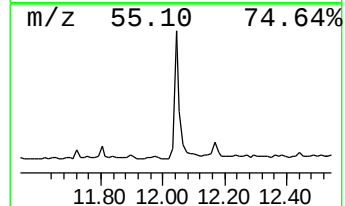
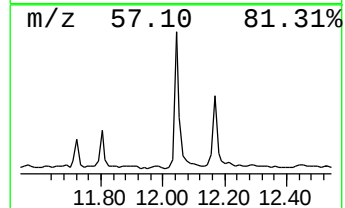
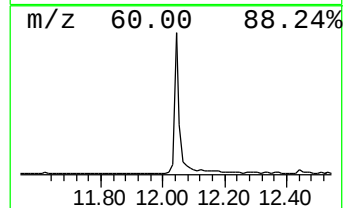
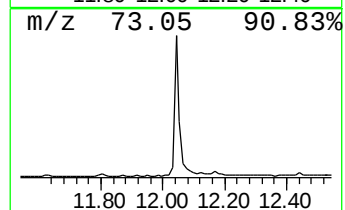
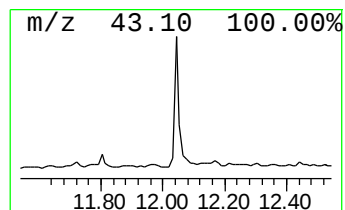
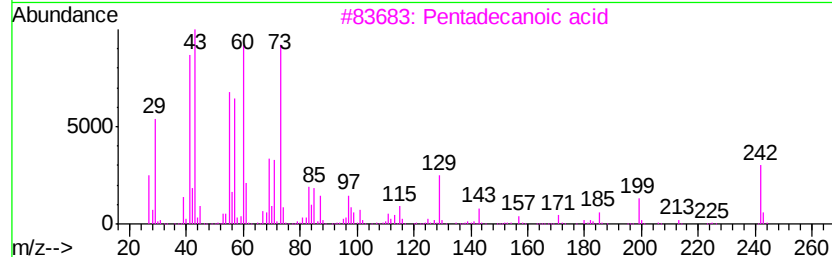
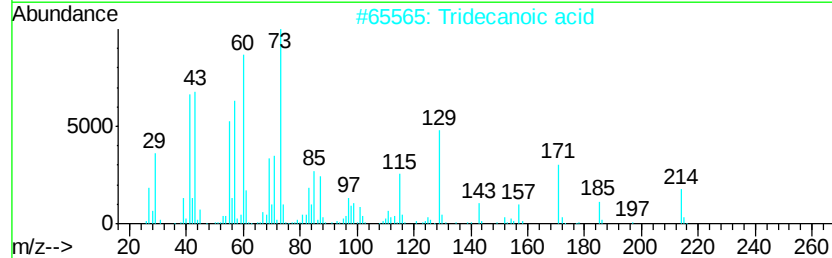
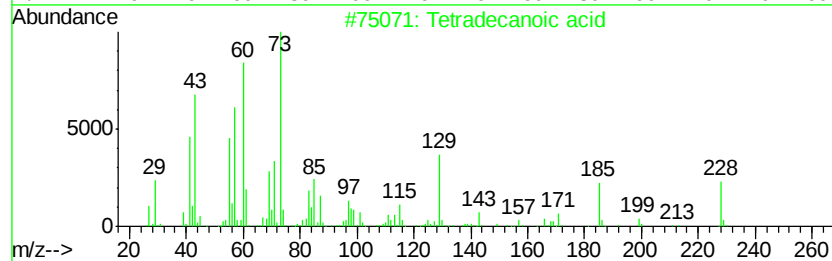
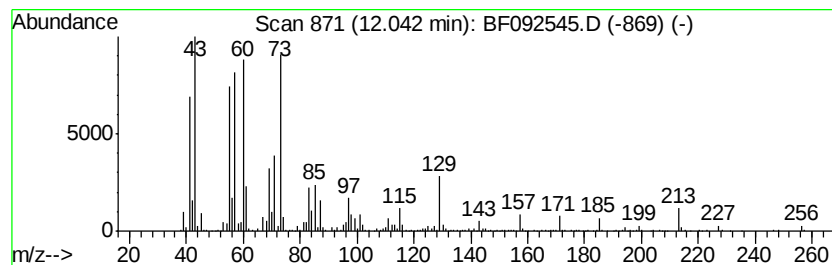
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Tetradecanoic acid Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.04 | 9.38 ng | 740111 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 94   |
| 2    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 93   |
| 3    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 91   |
| 4    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 91   |
| 5    |    | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 80   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

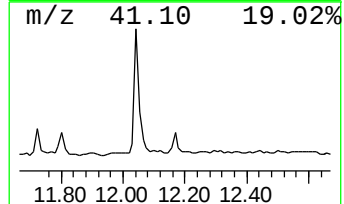
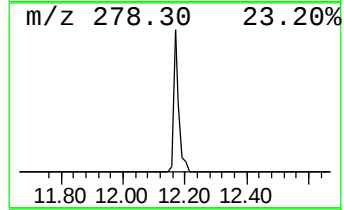
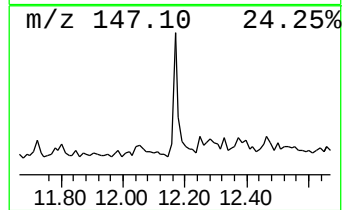
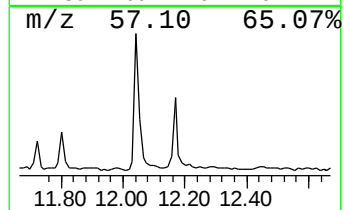
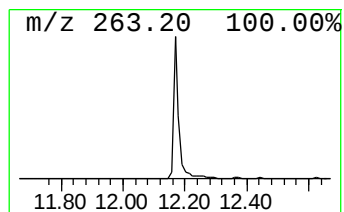
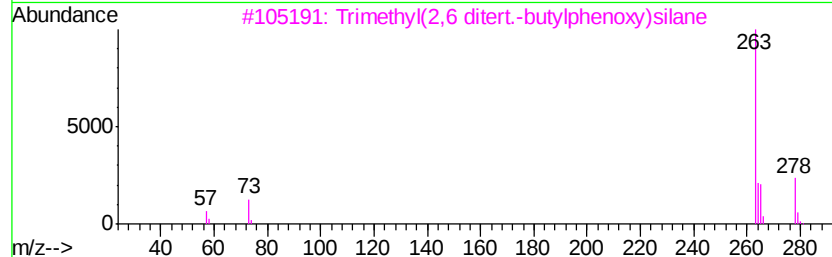
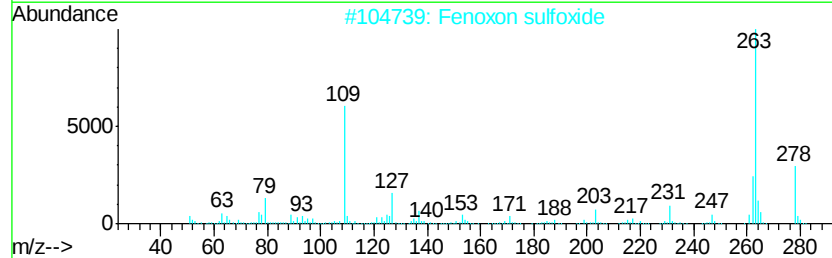
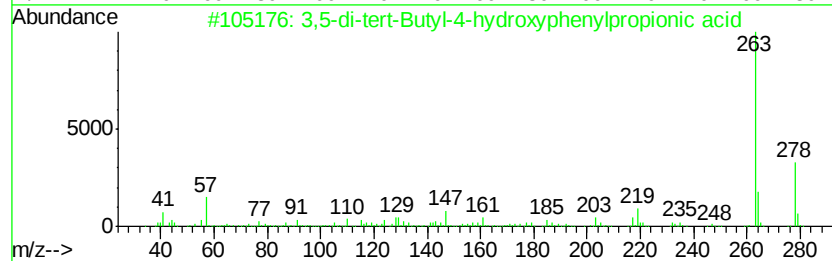
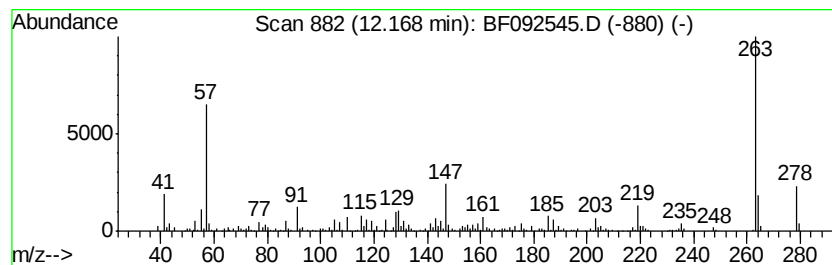
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.17 | 2.79 ng | 220097 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                         | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 3,5-di-tert-Butyl-4-hydroxyphenyl... | 278 | C17H26O3   | 020170-32-5  | 93   |
| 2    |    | Fenoxon sulfoxide                    | 278 | C10H15O5PS | 006552-13-2  | 50   |
| 3    |    | Trimethyl(2,6 di-tert.-butylpheno... | 278 | C17H30OSi  | 010416-73-6  | 47   |
| 4    |    | Ethanone, [1-(2,5-diphenyl-2H-1,...  | 263 | C16H13N3O  | 322647-34-7  | 43   |
| 5    |    | Iron, (.eta.-5-cyclopentadienyl)...  | 263 | C15H13FeN  | 1000164-28-2 | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

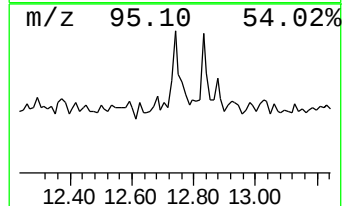
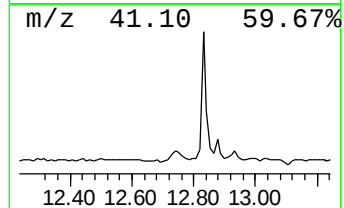
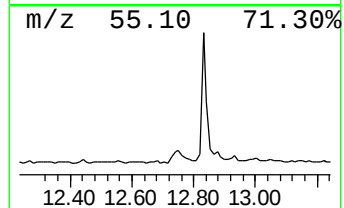
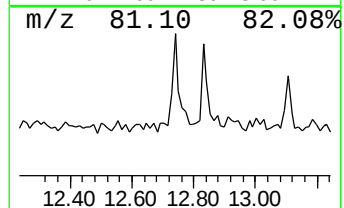
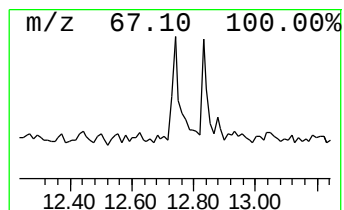
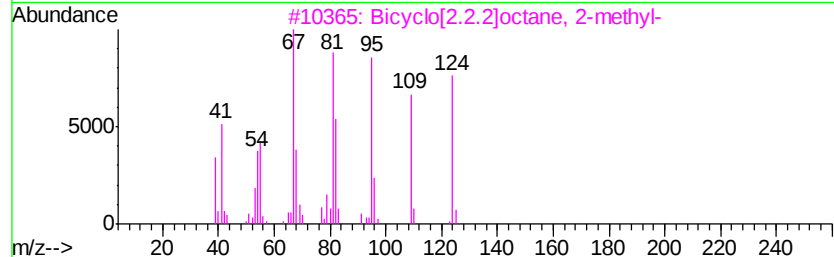
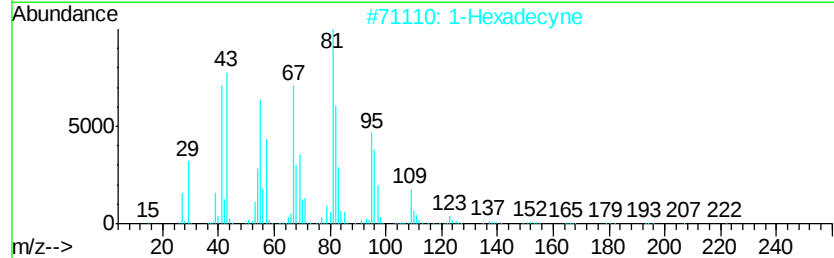
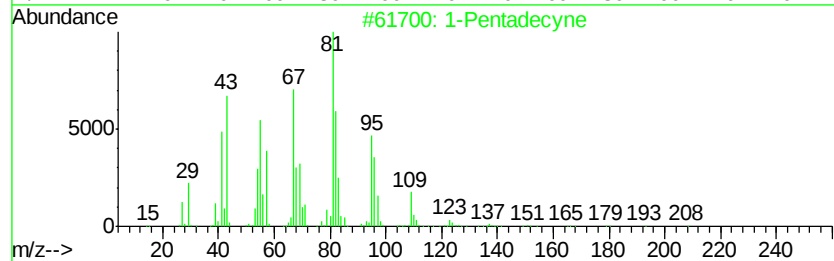
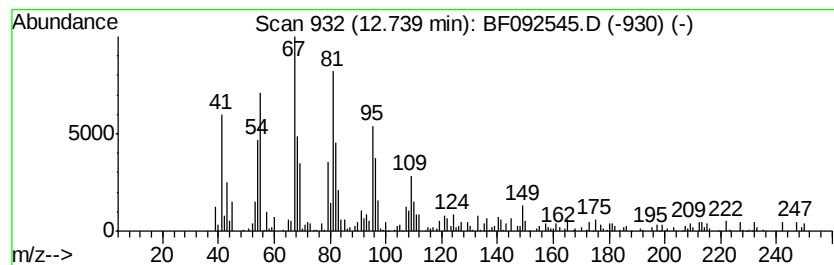
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 1-Pentadecyne Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.74 | 2.55 ng | 201179 | Phenanthrene-d10 | 11.52 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Pentadecyne                   | 208 | C15H28  | 000765-13-9 | 93   |
| 2    |    |   | 1-Hexadecyne                    | 222 | C16H30  | 000629-74-3 | 91   |
| 3    |    |   | Bicyclo[2.2.2]octane, 2-methyl- | 124 | C9H16   | 000766-53-0 | 91   |
| 4    |    |   | 1,9-Cyclohexadecadiene          | 220 | C16H28  | 004277-06-9 | 91   |
| 5    |    |   | 1-Tetradecyne                   | 194 | C14H26  | 000765-10-6 | 87   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

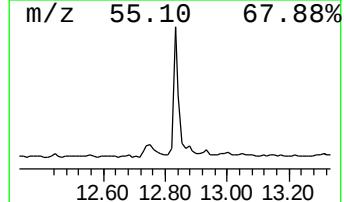
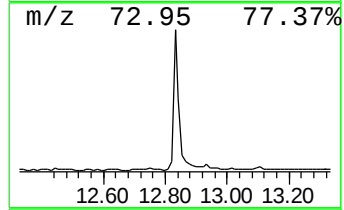
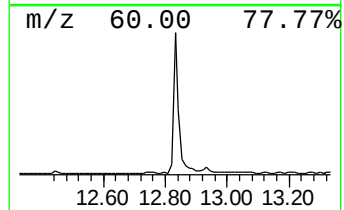
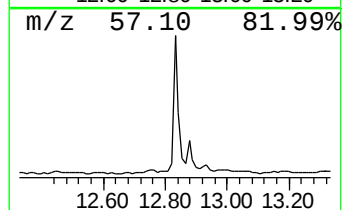
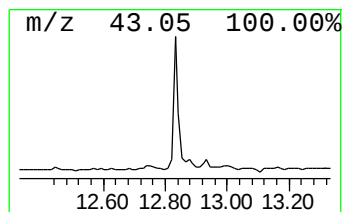
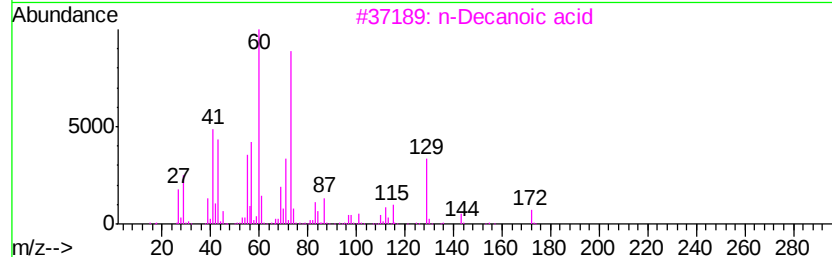
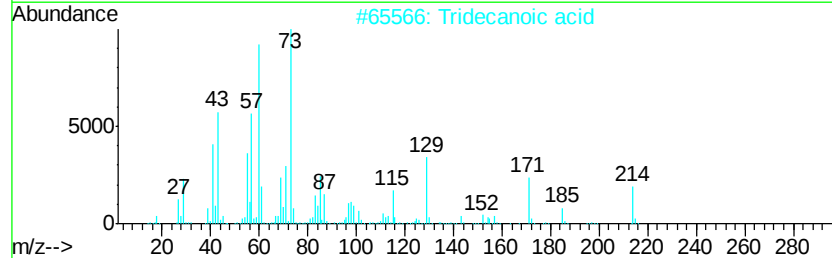
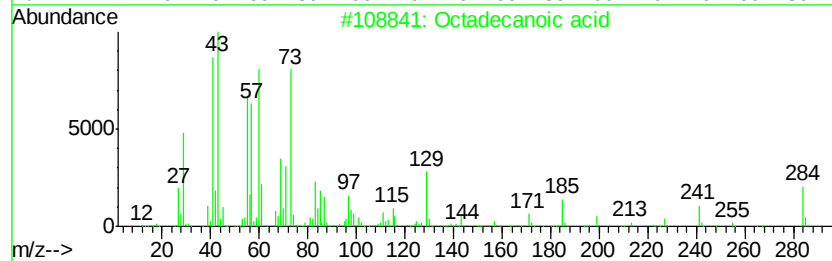
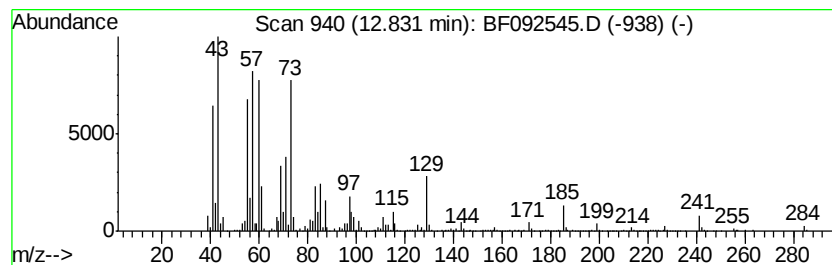
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Octadecanoic acid Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.83 | 13.48 ng | 1064140 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID          | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------|-----|----------|-------------|------|
| 1    | 5  | Octadecanoic acid     | 284 | C18H36O2 | 000057-11-4 | 94   |
| 2    |    | Tridecanoic acid      | 214 | C13H26O2 | 000638-53-9 | 81   |
| 3    |    | n-Decanoic acid       | 172 | C10H20O2 | 000334-48-5 | 74   |
| 4    |    | Pentadecanoic acid    | 242 | C15H30O2 | 001002-84-2 | 64   |
| 5    |    | Hexadecane, 1-chloro- | 260 | C16H33Cl | 004860-03-1 | 35   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

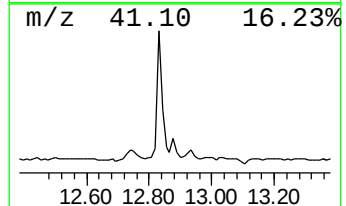
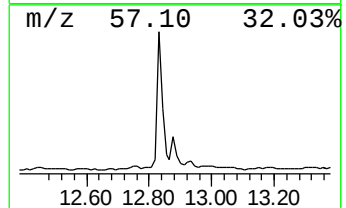
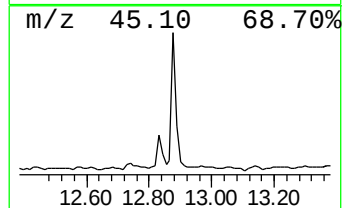
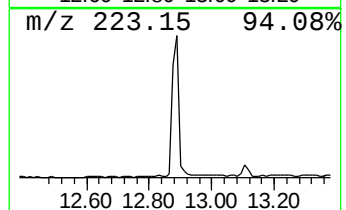
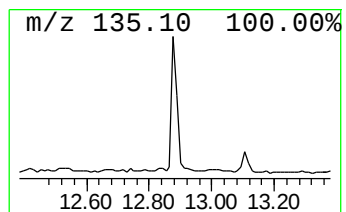
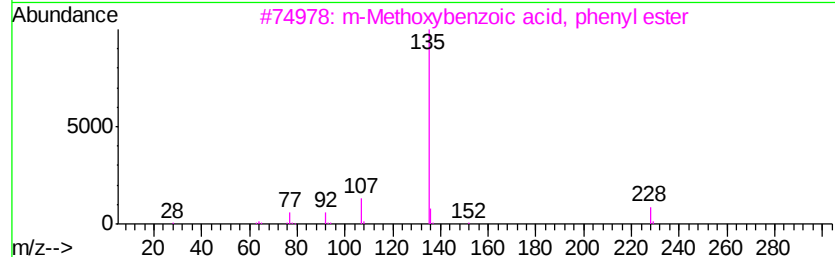
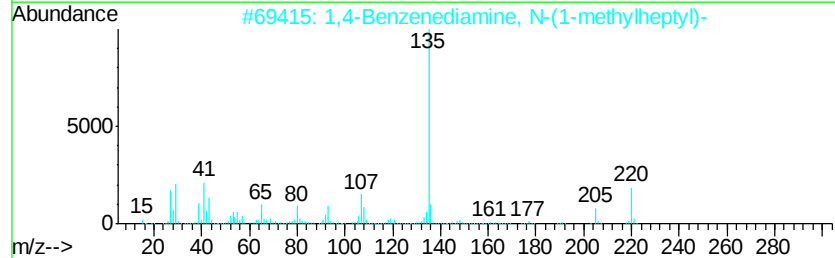
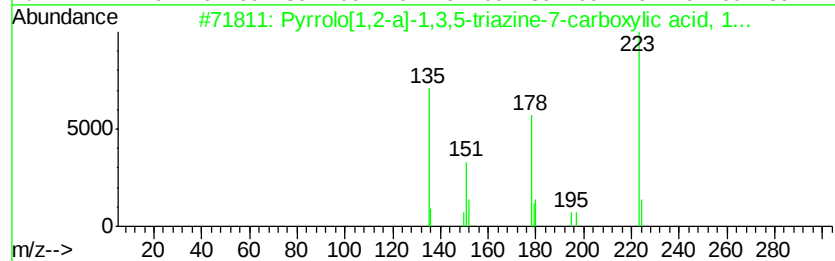
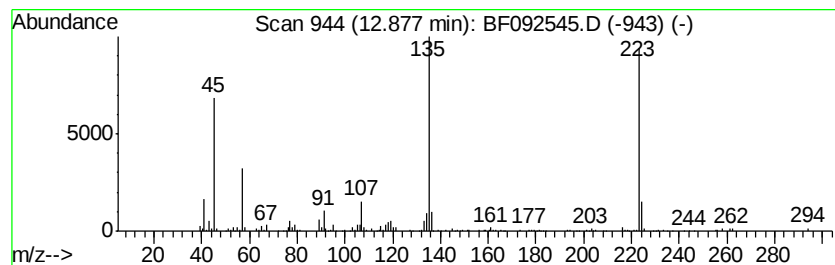
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Pyrrolo[1,2-a]-1,3,5-triazi... Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.88 | 4.02 ng | 286190 | Chrysene-d12     | 14.16 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Pyrrolo[1,2-a]-1,3,5-triazine-7-... | 223 | C9H9N3O4 | 054449-89-7 | 90   |
| 2    |    | 1,4-Benzenediamine, N-(1-methylh... | 220 | C14H24N2 | 039563-50-3 | 35   |
| 3    |    | m-Methoxybenzoic acid, phenyl ester | 228 | C14H12O3 | 065853-67-0 | 35   |
| 4    |    | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O  | 000080-46-6 | 27   |
| 5    |    | 2-Methoxycarbonyl-2-methylbrendane  | 194 | C12H18O2 | 148191-16-6 | 27   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

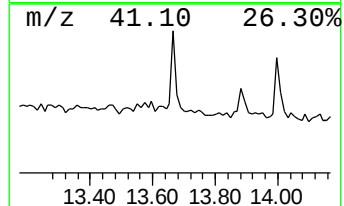
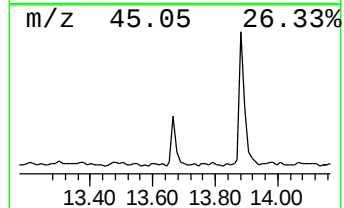
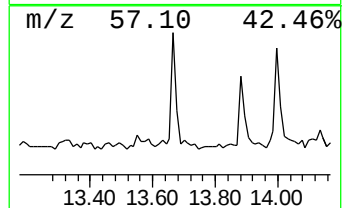
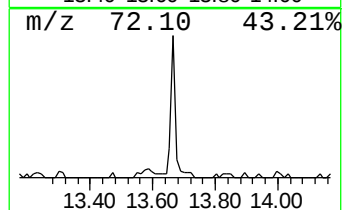
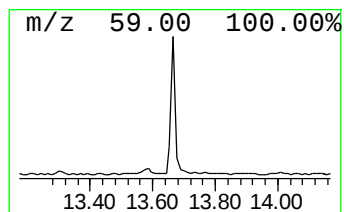
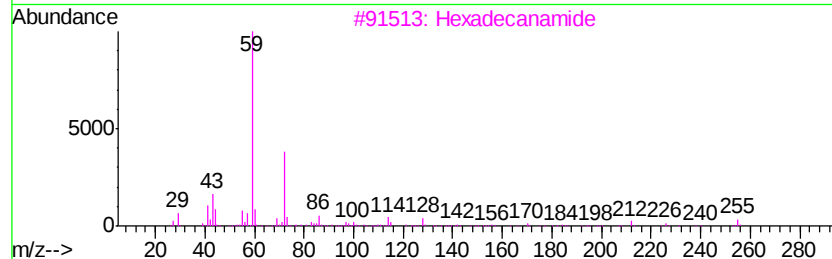
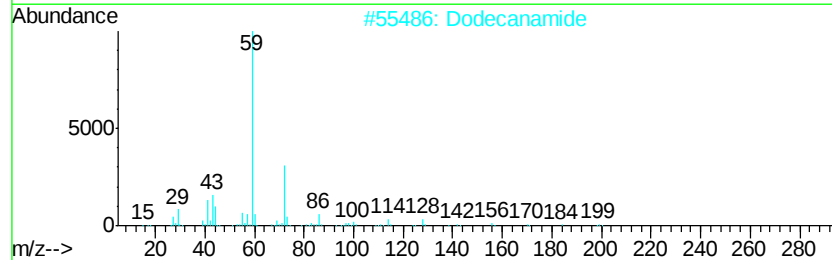
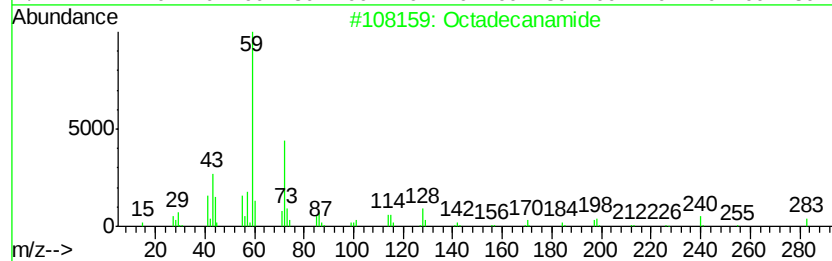
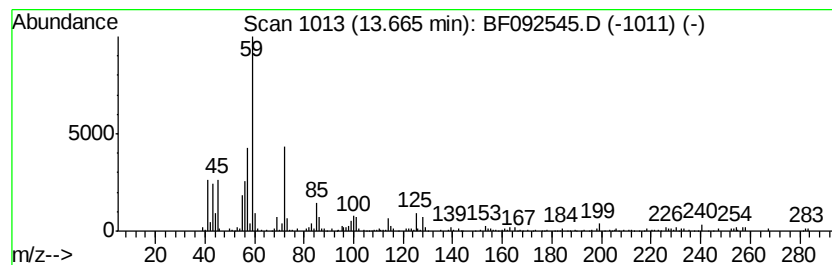
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Octadecanamide Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.67 | 2.07 ng | 147418 | Chrysene-d12     | 14.16 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Octadecanamide                      | 283 | C18H37NO   | 000124-26-5 | 70   |
| 2    |    | Dodecanamide                        | 199 | C12H25NO   | 001120-16-7 | 64   |
| 3    |    | Hexadecanamide                      | 255 | C16H33NO   | 000629-54-9 | 64   |
| 4    |    | Benzeneethanamine, 2-fluoro-.bet... | 229 | C11H16FN03 | 061338-98-5 | 53   |
| 5    |    | Tetradecanamide                     | 227 | C14H29NO   | 000638-58-4 | 50   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\  
Data File : BF092545.D  
Acq On : 27 Jan 2017 00:37  
Operator : SJ/MA  
Sample : I1451-04  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
W-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

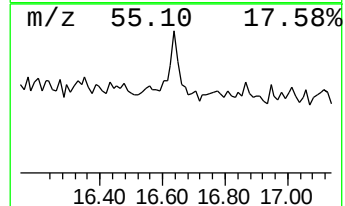
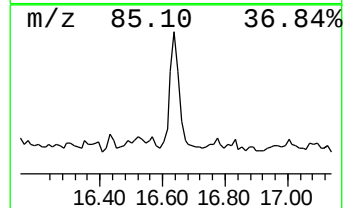
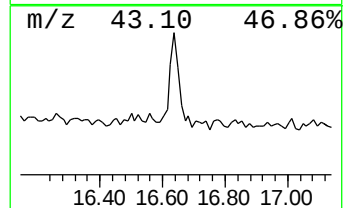
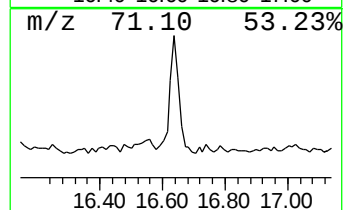
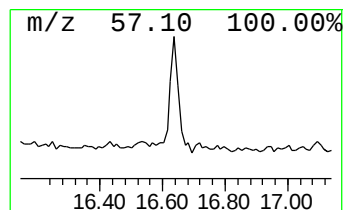
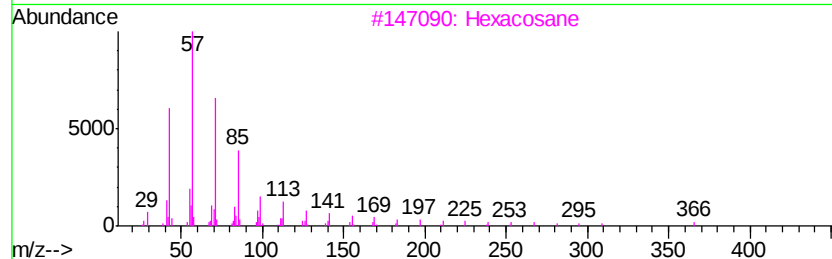
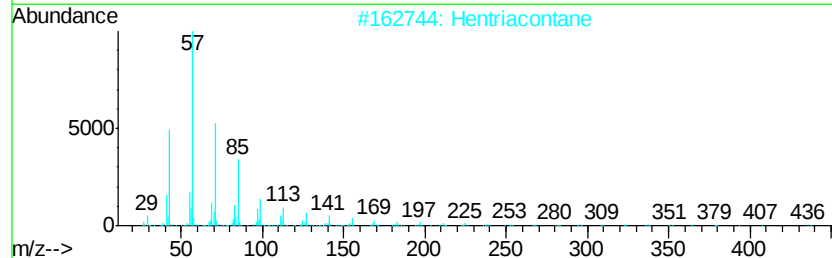
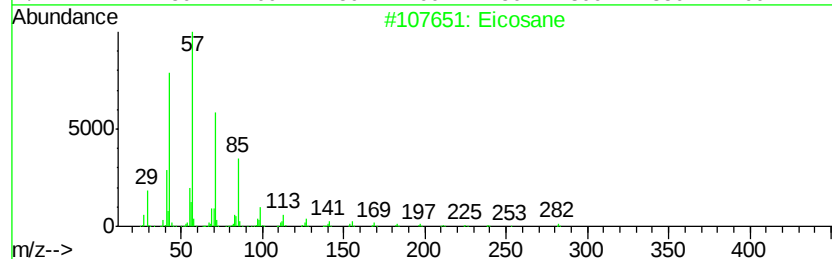
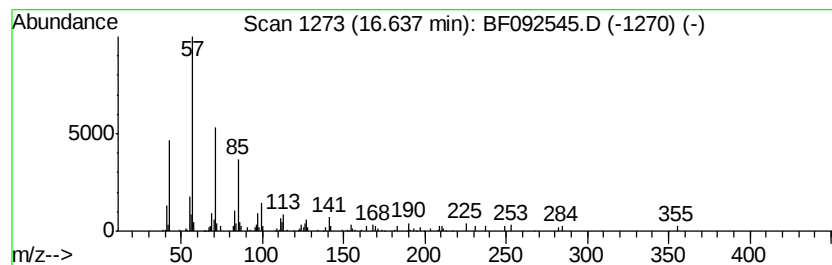
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Eicosane Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.64 | 2.66 ng | 142368 | Perylene-d12     | 15.64 |

| Hit# | of | Tentative ID   | MW  | MolForm | CAS#        | Qual |
|------|----|----------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane       | 282 | C20H42  | 000112-95-8 | 83   |
| 2    |    | Hentriacontane | 437 | C31H64  | 000630-04-6 | 80   |
| 3    |    | Hexacosane     | 366 | C26H54  | 000630-01-3 | 72   |
| 4    |    | Heptacosane    | 380 | C27H56  | 000593-49-7 | 72   |
| 5    |    | Hexadecane     | 226 | C16H34  | 000544-76-3 | 72   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF012617\  
 Data File : BF092545.D  
 Acq On : 27 Jan 2017 00:37  
 Operator : SJ/MA  
 Sample : I1451-04  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 W-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF012417.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy... | 5.15  | 12.3    | ng    | 628088   | 1                     | 6.97  | 1021510 | 20.0 |
| unknown6.73          | 6.73  | 76.2    | ng    | 3894040  | 1                     | 6.97  | 1021510 | 20.0 |
| Acetamide, N-(2,4... | 10.95 | 4.6     | ng    | 361198   | 4                     | 11.52 | 1578470 | 20.0 |
| Phenol, 4-(1,1,3,... | 10.99 | 7.1     | ng    | 562163   | 4                     | 11.52 | 1578470 | 20.0 |
| unknown11.06         | 11.06 | 11.4    | ng    | 902228   | 4                     | 11.52 | 1578470 | 20.0 |
| unknown11.13         | 11.13 | 7.8     | ng    | 614707   | 4                     | 11.52 | 1578470 | 20.0 |
| 2-Methyl-4-hydrox... | 11.17 | 8.0     | ng    | 629413   | 4                     | 11.52 | 1578470 | 20.0 |
| Phenol, 4-(1,1-di... | 11.21 | 9.5     | ng    | 749247   | 4                     | 11.52 | 1578470 | 20.0 |
| unknown11.25         | 11.25 | 4.1     | ng    | 321720   | 4                     | 11.52 | 1578470 | 20.0 |
| Ethanol, 2-[4-(1,... | 11.72 | 2.3     | ng    | 177960   | 4                     | 11.52 | 1578470 | 20.0 |
| Benzoic acid, 3,5... | 11.80 | 2.7     | ng    | 210909   | 4                     | 11.52 | 1578470 | 20.0 |
| Tetradecanoic acid   | 12.04 | 9.4     | ng    | 740111   | 4                     | 11.52 | 1578470 | 20.0 |
| 3,5-di-tert-Butyl... | 12.17 | 2.8     | ng    | 220097   | 4                     | 11.52 | 1578470 | 20.0 |
| 1-Pentadecyne        | 12.74 | 2.5     | ng    | 201179   | 4                     | 11.52 | 1578470 | 20.0 |
| Octadecanoic acid    | 12.83 | 13.5    | ng    | 1064140  | 4                     | 11.52 | 1578470 | 20.0 |
| Pyrrolo[1,2-a]-1,... | 12.88 | 4.0     | ng    | 286190   | 5                     | 14.16 | 1424950 | 20.0 |
| Octadecanamide       | 13.67 | 2.1     | ng    | 147418   | 5                     | 14.16 | 1424950 | 20.0 |
| Eicosane             | 16.64 | 2.7     | ng    | 142368   | 6                     | 15.64 | 1072270 | 20.0 |