

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\  
 Data File : BF093203.D  
 Acq On : 27 Feb 2017 12:37  
 Operator : SJ/MA  
 Sample : I1954-09  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CLP-B-01(0-5)

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-BF013017.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         | 4.990       | 251           | 254         | 262          | rBV      | 1365388        | 1930376       | 43.22%          | 4.720%        |
| 2         | 5.424       | 290           | 292         | 295          | rBV      | 3496745        | 4030381       | 90.23%          | 9.854%        |
| 3         | 6.476       | 381           | 384         | 387          | rBV      | 3407415        | 4248339       | 95.11%          | 10.387%       |
| 4         | 6.602       | 392           | 395         | 397          | rBV      | 4061523        | 4420346       | 98.97%          | 10.808%       |
| 5         | 6.830       | 412           | 415         | 417          | rBV      | 905713         | 1020734       | 22.85%          | 2.496%        |
| 6         | 6.979       | 426           | 428         | 431          | rBV      | 2367591        | 3309207       | 74.09%          | 8.091%        |
| 7         | 7.402       | 462           | 465         | 467          | rBV      | 2878745        | 2873662       | 64.34%          | 7.026%        |
| 8         | 8.122       | 525           | 528         | 530          | rBV      | 1517595        | 1410221       | 31.57%          | 3.448%        |
| 9         | 9.208       | 620           | 623         | 625          | rBV      | 4537026        | 4466545       | 100.00%         | 10.921%       |
| 10        | 9.608       | 655           | 658         | 659          | rBV      | 236290         | 295032        | 6.61%           | 0.721%        |
| 11        | 9.813       | 675           | 676         | 680          | rBV3     | 56677          | 82198         | 1.84%           | 0.201%        |
| 12        | 9.882       | 680           | 682         | 684          | rVB      | 1582709        | 1373040       | 30.74%          | 3.357%        |
| 13        | 10.316      | 716           | 720         | 721          | rBV2     | 125712         | 154691        | 3.46%           | 0.378%        |
| 14        | 10.533      | 736           | 739         | 742          | rBV      | 34900          | 52300         | 1.17%           | 0.128%        |
| 15        | 10.682      | 749           | 752         | 754          | rBV      | 1742549        | 2049690       | 45.89%          | 5.012%        |
| 16        | 10.785      | 760           | 761         | 766          | rVB      | 204534         | 303195        | 6.79%           | 0.741%        |
| 17        | 11.231      | 799           | 800         | 802          | rBV      | 94409          | 139785        | 3.13%           | 0.342%        |
| 18        | 11.368      | 810           | 812         | 816          | rBV      | 1260807        | 1242166       | 27.81%          | 3.037%        |
| 19        | 11.608      | 831           | 833         | 835          | rBV3     | 35685          | 65278         | 1.46%           | 0.160%        |
| 20        | 11.665      | 835           | 838         | 839          | rVV      | 87879          | 98926         | 2.21%           | 0.242%        |
| 21        | 11.871      | 854           | 856         | 857          | rBV      | 46487          | 52567         | 1.18%           | 0.129%        |
| 22        | 11.974      | 864           | 865         | 869          | rVB      | 43567          | 92064         | 2.06%           | 0.225%        |
| 23        | 12.065      | 871           | 873         | 875          | rBV      | 61778          | 67282         | 1.51%           | 0.165%        |
| 24        | 12.168      | 878           | 882         | 883          | rBV2     | 25994          | 53005         | 1.19%           | 0.130%        |
| 25        | 12.294      | 891           | 893         | 896          | rBV2     | 31238          | 44811         | 1.00%           | 0.110%        |
| 26        | 12.419      | 902           | 904         | 905          | rBV      | 95903          | 108180        | 2.42%           | 0.265%        |
| 27        | 12.454      | 905           | 907         | 910          | rVB2     | 81418          | 115678        | 2.59%           | 0.283%        |
| 28        | 12.511      | 910           | 912         | 913          | rBV      | 35249          | 54705         | 1.22%           | 0.134%        |
| 29        | 12.579      | 916           | 918         | 920          | rVB      | 224712         | 192365        | 4.31%           | 0.470%        |
| 30        | 12.774      | 933           | 935         | 936          | rBV      | 155612         | 162753        | 3.64%           | 0.398%        |
| 31        | 12.808      | 936           | 938         | 941          | rVV      | 282463         | 334031        | 7.48%           | 0.817%        |
| 32        | 12.957      | 948           | 951         | 953          | rBV      | 2579493        | 3033738       | 67.92%          | 7.418%        |
| 33        | 13.128      | 963           | 966         | 968          | rBV3     | 83507          | 155693        | 3.49%           | 0.381%        |
| 34        | 13.174      | 968           | 970         | 974          | rBV3     | 62787          | 151471        | 3.39%           | 0.370%        |

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ClientSampleId :  
CLP-B-01(0-5)

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-BF013017.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 13.471 | 995  | 996  | 998  | rBV  | 78762  | 89475   | 2.00%  | 0.219% |
| 36 | 13.517 | 998  | 1000 | 1002 | rBV2 | 84664  | 113360  | 2.54%  | 0.277% |
| 37 | 13.848 | 1027 | 1029 | 1033 | rVB4 | 241212 | 369963  | 8.28%  | 0.905% |
| 38 | 14.008 | 1040 | 1043 | 1044 | rBV  | 835522 | 1157772 | 25.92% | 2.831% |
| 39 | 15.437 | 1166 | 1168 | 1171 | rBV  | 679125 | 984096  | 22.03% | 2.406% |

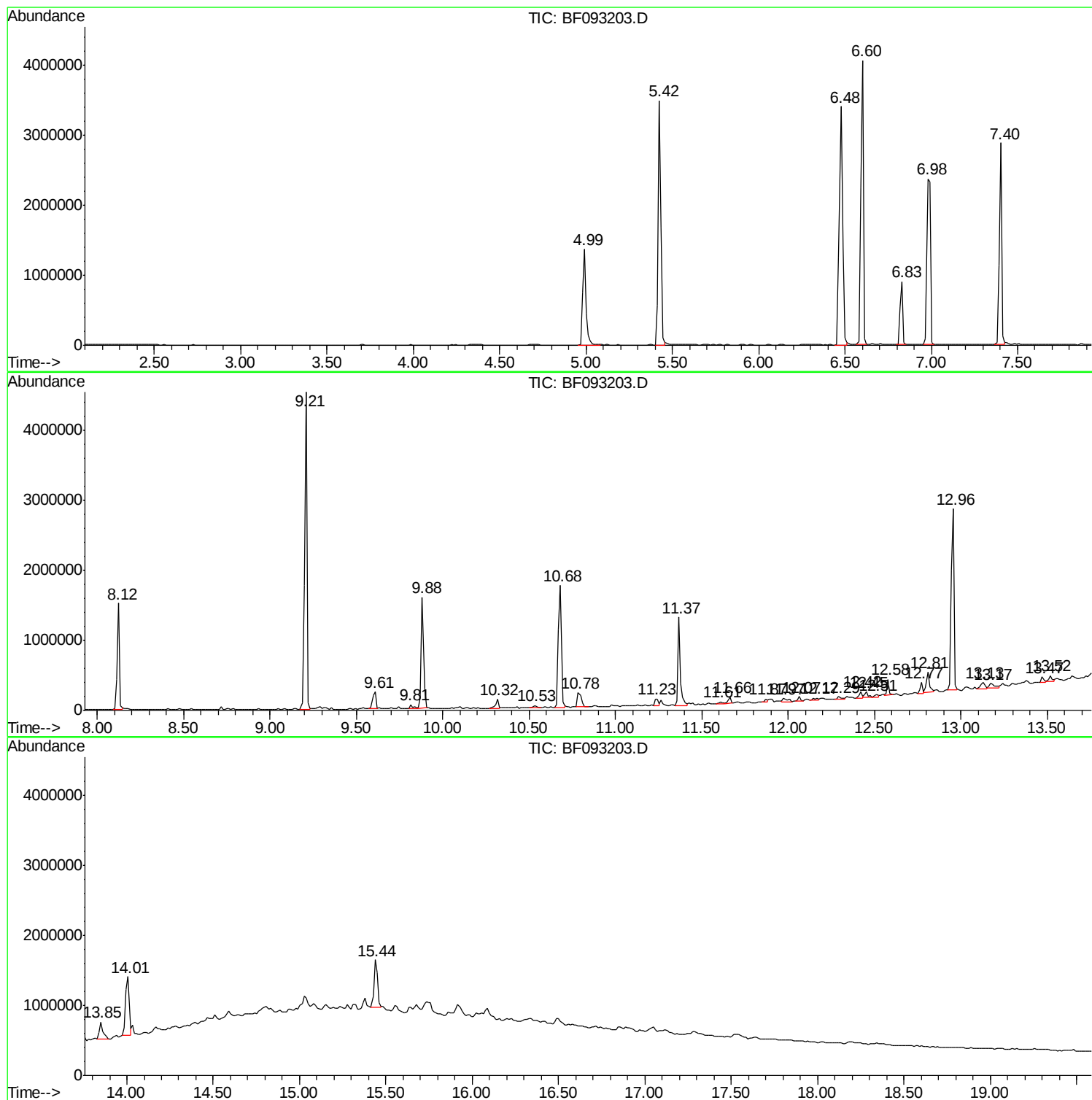
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ClientSampled :  
CLP-B-01(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF013017.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-BF013017.M  
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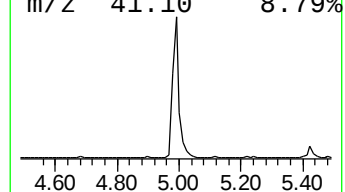
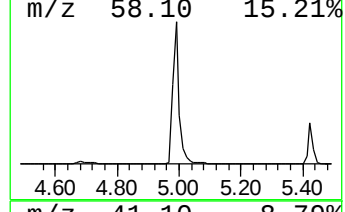
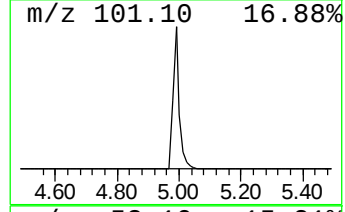
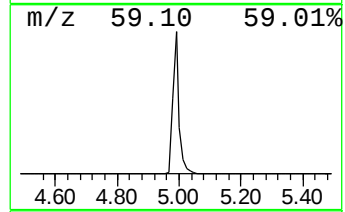
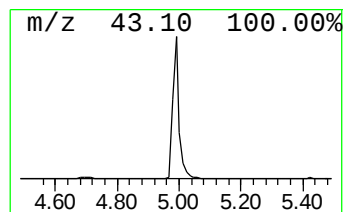
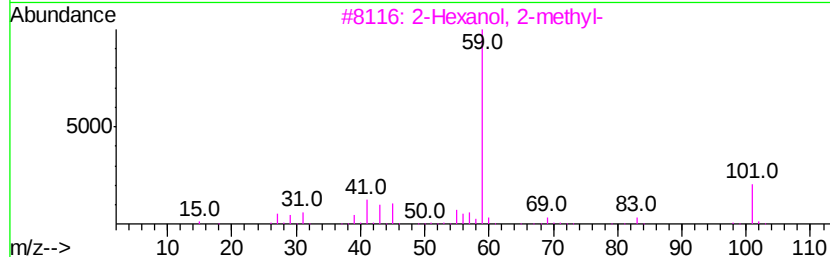
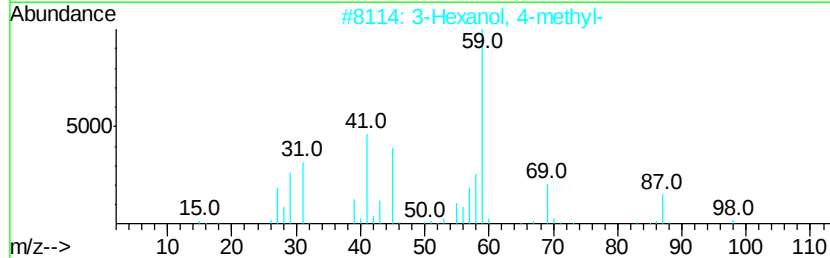
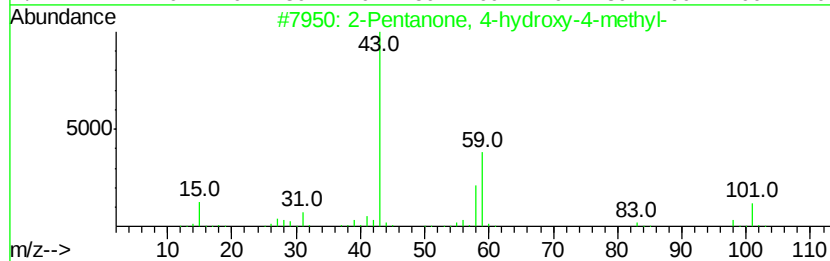
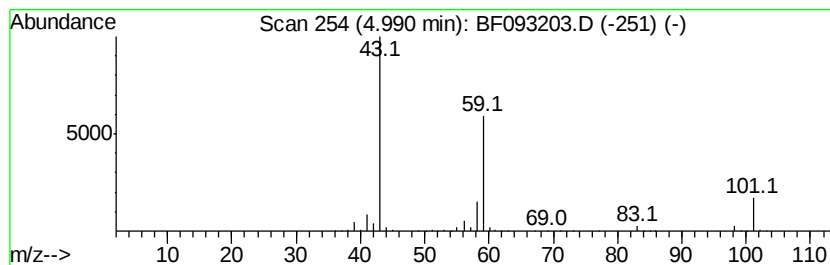
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 4.99 | 37.82 ng | 1930380 | 1,4-Dichlorobenzene-d4 | 6.83 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 59   |
| 2    |    | 3-Hexanol, 4-methyl-                 | 116 | C7H16O   | 000615-29-2 | 33   |
| 3    |    | 2-Hexanol, 2-methyl-                 | 116 | C7H16O   | 000625-23-0 | 33   |
| 4    |    | Propane, 2-methyl-2-(1-methylethyl)- | 116 | C7H16O   | 017348-59-3 | 33   |
| 5    |    | Acetic acid, cyano-, 1,1-dimethyl-   | 141 | C7H11NO2 | 001116-98-9 | 23   |



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CLP-B-01(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

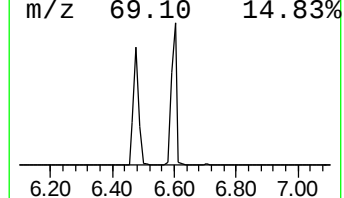
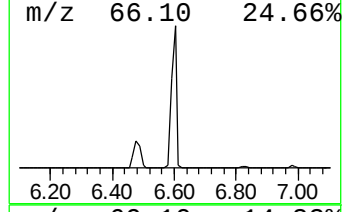
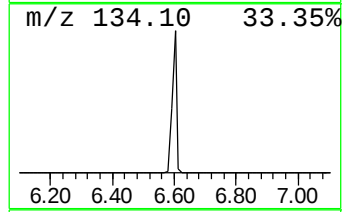
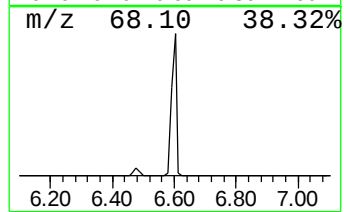
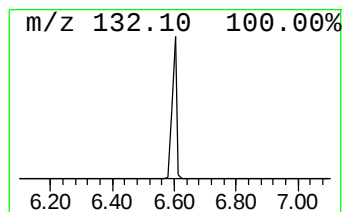
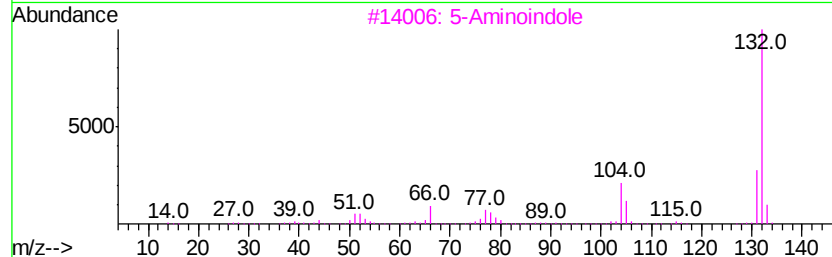
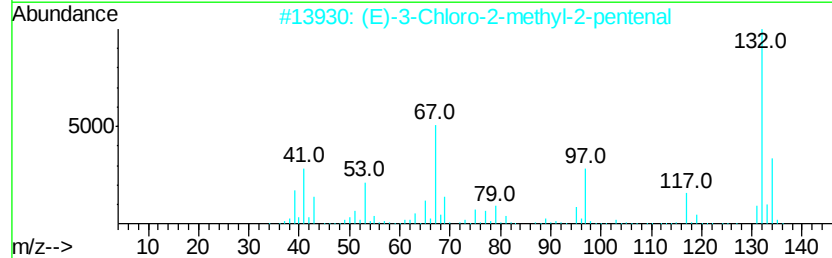
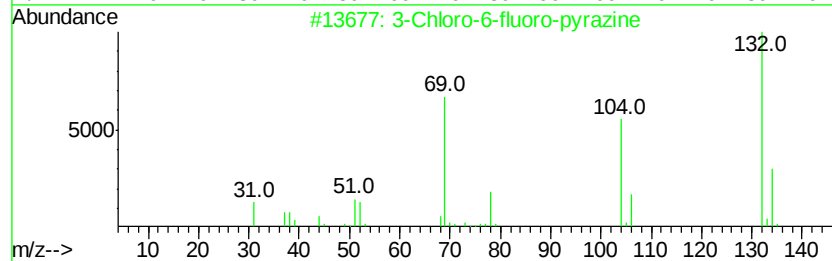
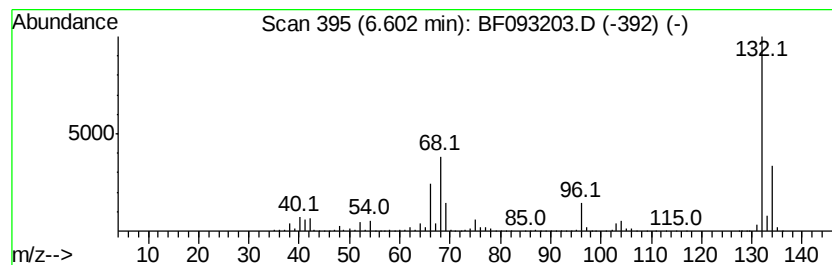
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown6.60 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.60 | 86.61 ng | 4420350 | 1,4-Dichlorobenzene-d4 | 6.83 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | Benzene, 1,3,5-trifluoro-        | 132 | C6H3F3    | 000372-38-3  | 9    |



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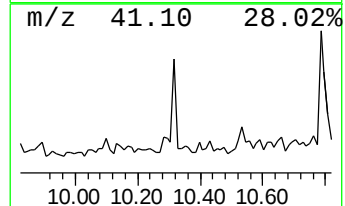
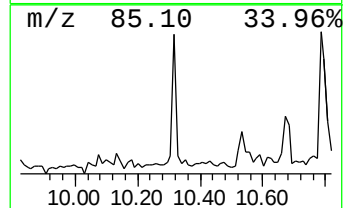
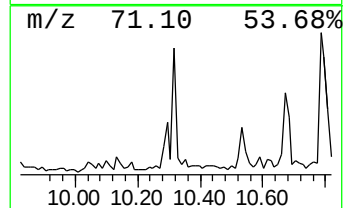
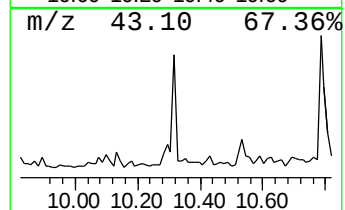
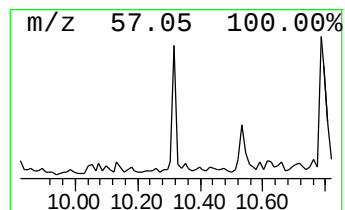
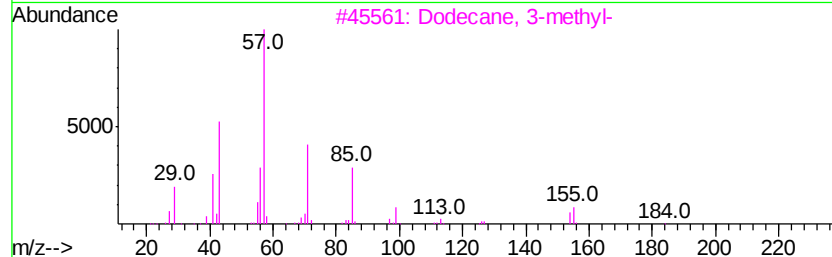
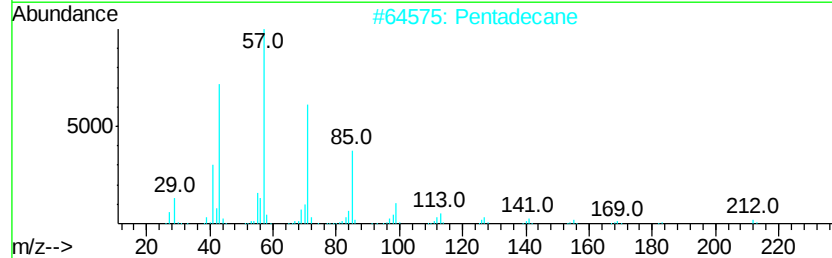
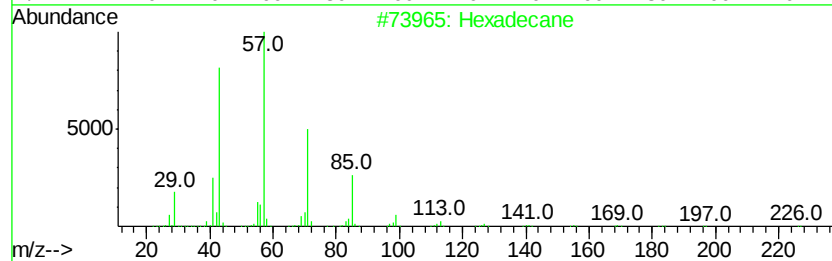
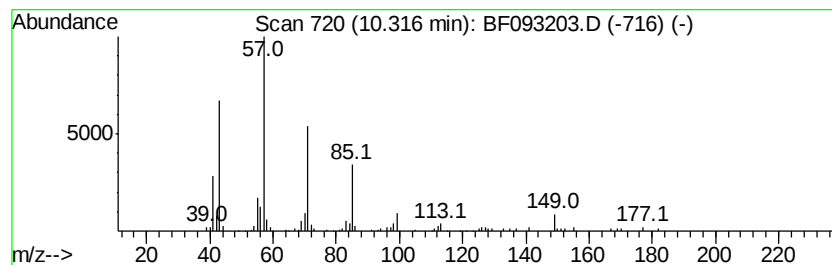
Instrument :  
BNA\_F  
ClientSampleId :  
CLP-B-01(0-5)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Hexadecane Concentration Rank 8

| R.T.    | EstConc                  | Area         | Relative to ISTD | R.T.      |
|---------|--------------------------|--------------|------------------|-----------|
| 10.32   | 2.25 ng                  | 154691       | Acenaphthene-d10 | 9.88      |
| Hit# of | 5                        | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Hexadecane               | 226 C16H34   | 000544-76-3      | 86        |
| 2       | Pentadecane              | 212 C15H32   | 000629-62-9      | 80        |
| 3       | Dodecane, 3-methyl-      | 184 C13H28   | 017312-57-1      | 78        |
| 4       | Decane, 2,3,5-trimethyl- | 184 C13H28   | 062238-11-3      | 78        |
| 5       | Tridecane                | 184 C13H28   | 000629-50-5      | 78        |



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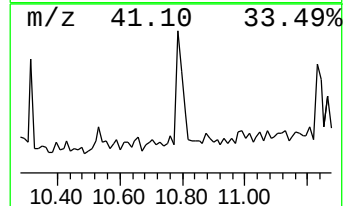
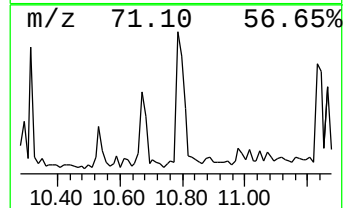
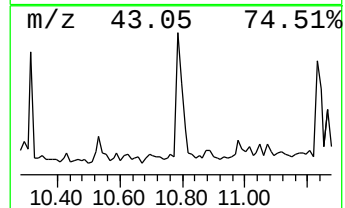
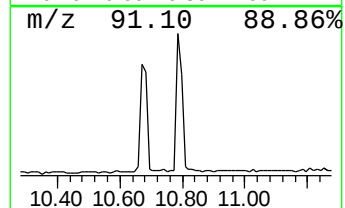
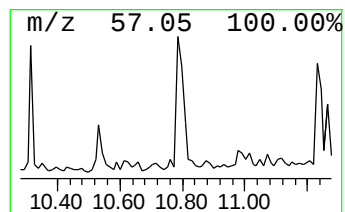
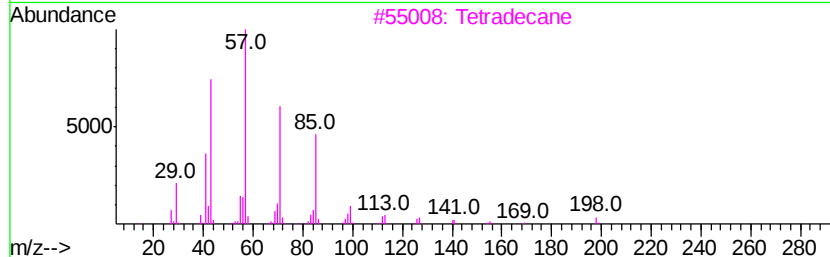
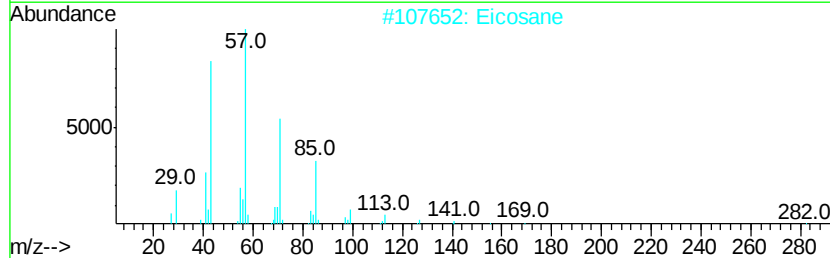
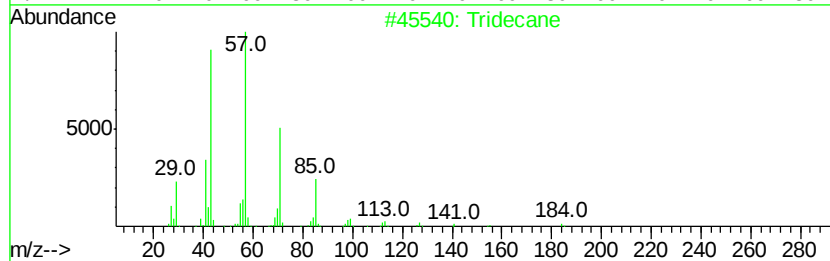
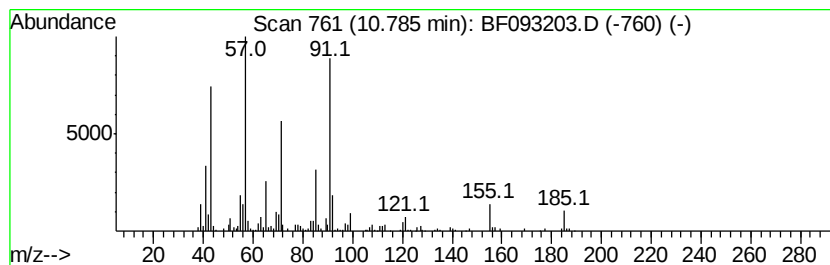
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Tridecane Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.78 | 4.88 ng | 303195 | Phenanthrene-d10 | 11.37 |

| Hit# of | 5           | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------|--------------|-----|---------|-------------|------|
| 1       | Tridecane   |              | 184 | C13H28  | 000629-50-5 | 50   |
| 2       | Eicosane    |              | 282 | C20H42  | 000112-95-8 | 49   |
| 3       | Tetradecane |              | 198 | C14H30  | 000629-59-4 | 46   |
| 4       | Hexadecane  |              | 226 | C16H34  | 000544-76-3 | 46   |
| 5       | Pentadecane |              | 212 | C15H32  | 000629-62-9 | 46   |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF013017.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

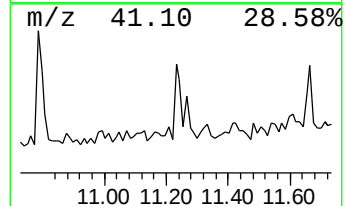
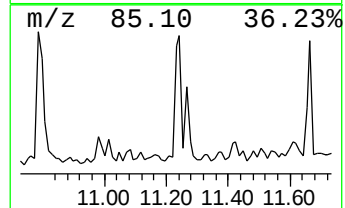
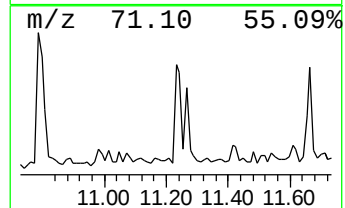
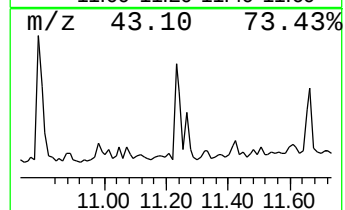
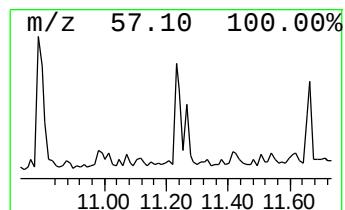
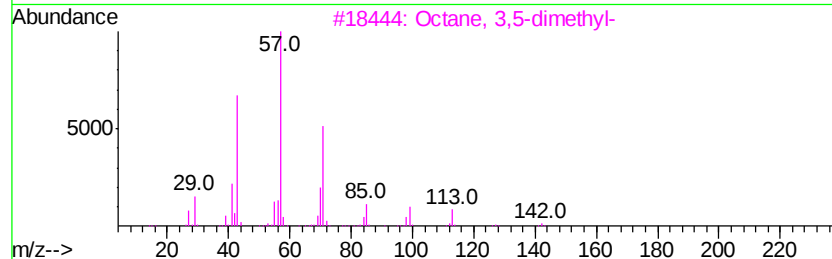
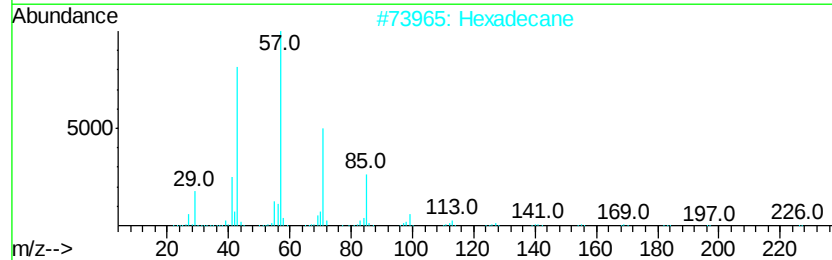
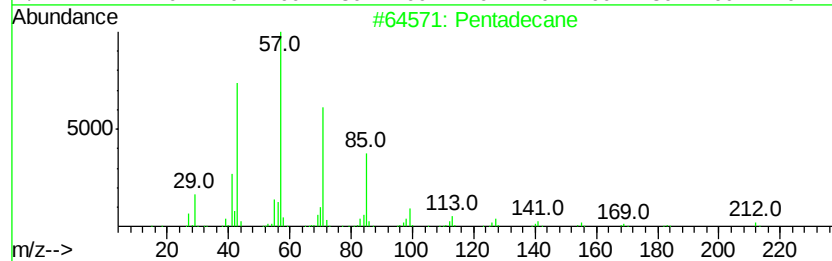
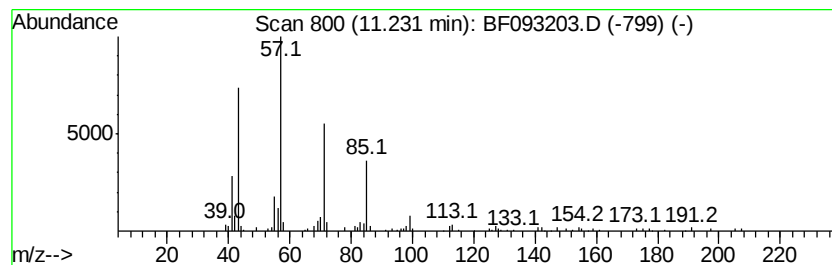
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Pentadecane Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.23 | 2.25 ng | 139785 | Phenanthrene-d10 | 11.37 |

| Hit# of | 5                      | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------------------|--------------|-----|---------|-------------|------|
| 1       | Pentadecane            |              | 212 | C15H32  | 000629-62-9 | 90   |
| 2       | Hexadecane             |              | 226 | C16H34  | 000544-76-3 | 90   |
| 3       | Octane, 3,5-dimethyl-  |              | 142 | C10H22  | 015869-93-9 | 83   |
| 4       | Tetradecane, 1-iodo-   |              | 324 | C14H29I | 019218-94-1 | 72   |
| 5       | 3-Heptanone, 4-methyl- |              | 128 | C8H16O  | 006137-11-7 | 70   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\  
Data File : BF093203.D  
Acq On : 27 Feb 2017 12:37  
Operator : SJ/MA  
Sample : I1954-09  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

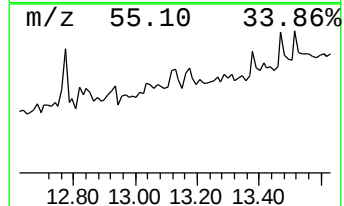
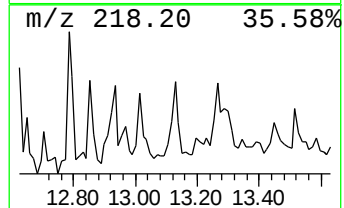
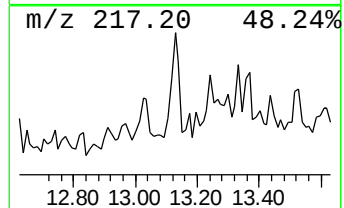
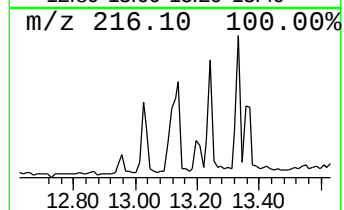
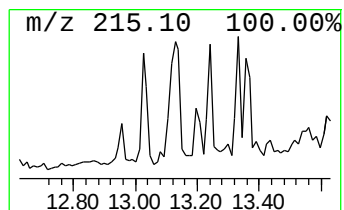
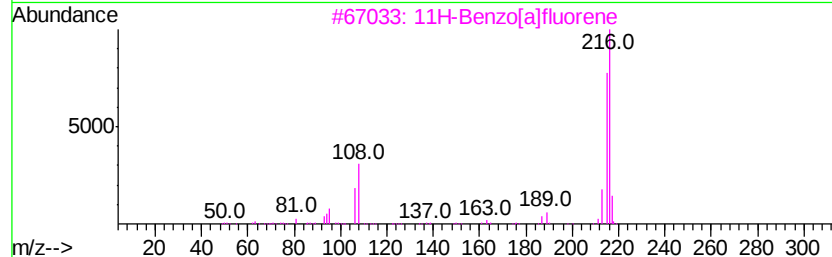
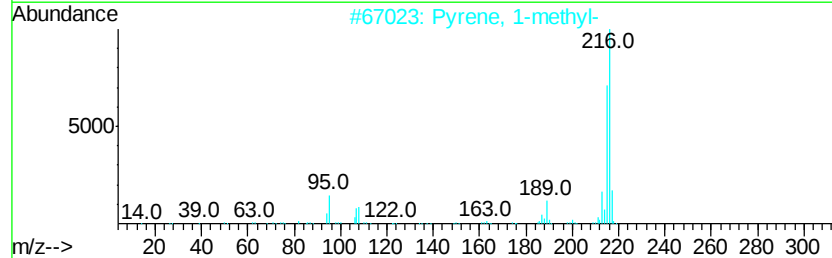
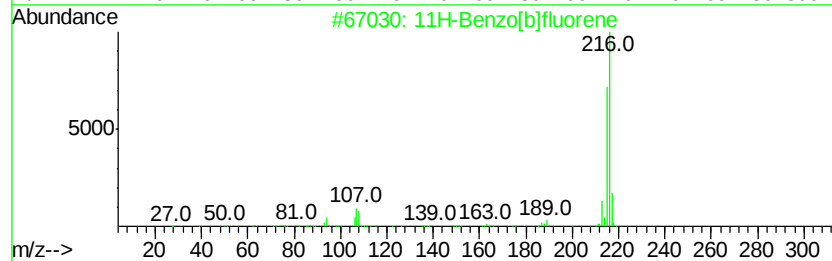
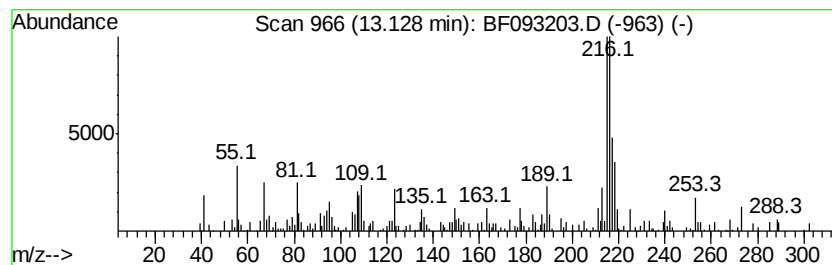
Instrument :  
BNA\_F  
ClientSampleId :  
CLP-B-01(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.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 11H-Benzo[b]fluorene Concentration Rank 6

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.     |             |      |
|---------|---------|-------------------------------------|------------------|----------|-------------|------|
| 13.13   | 2.69 ng | 155693                              | Chrysene-d12     | 14.01    |             |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm  | CAS#        | Qual |
| 1       |         | 11H-Benzo[blfluorene                | 216              | C17H12   | 000243-17-4 | 78   |
| 2       |         | Pyrene, 1-methyl-                   | 216              | C17H12   | 002381-21-7 | 60   |
| 3       |         | 11H-Benzo[alfluorene                | 216              | C17H12   | 000238-84-6 | 47   |
| 4       |         | Pyrene, 2-methyl-                   | 216              | C17H12   | 003442-78-2 | 41   |
| 5       |         | 1,3,5-Triazine-2,4,6-triamine, N... | 216              | C10H12N6 | 046731-79-7 | 38   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\  
Data File : BF093203.D  
Acq On : 27 Feb 2017 12:37  
Operator : SJ/MA  
Sample : I1954-09  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

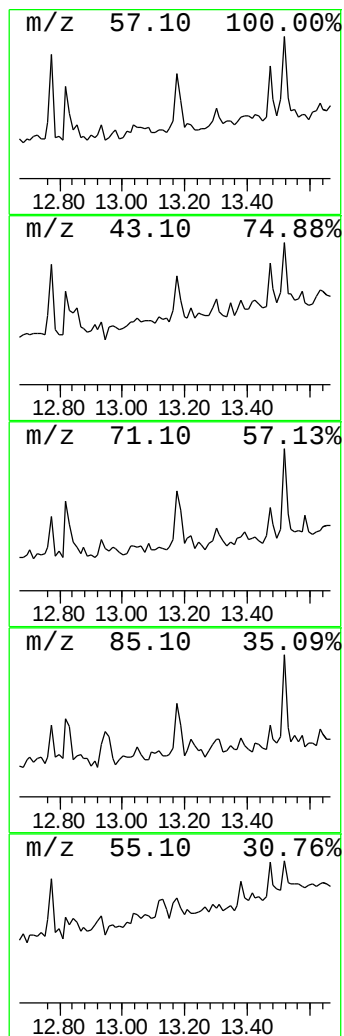
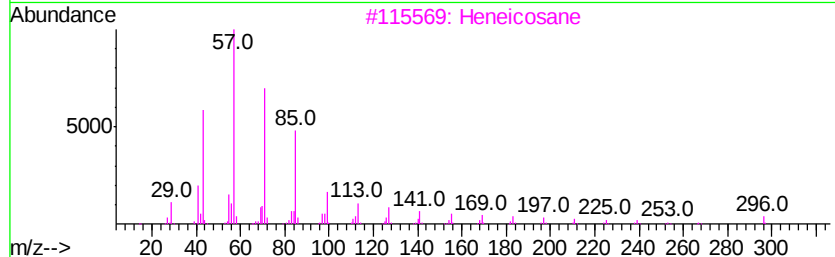
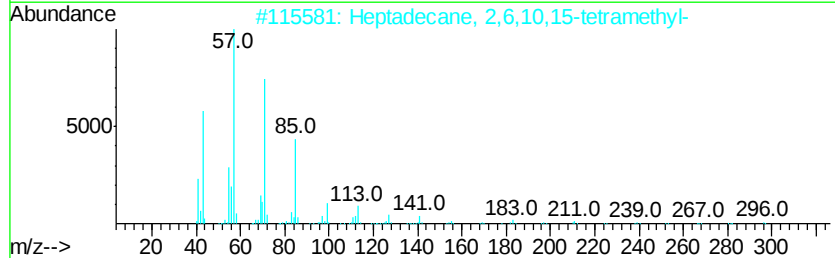
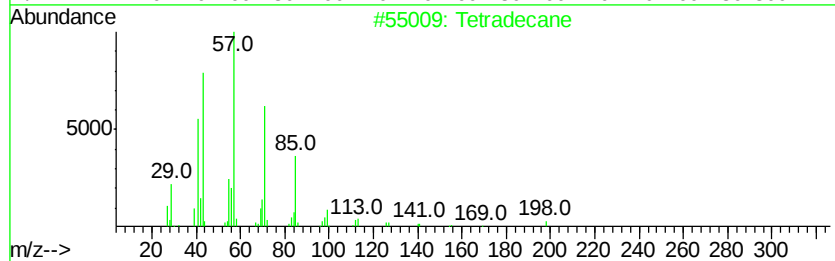
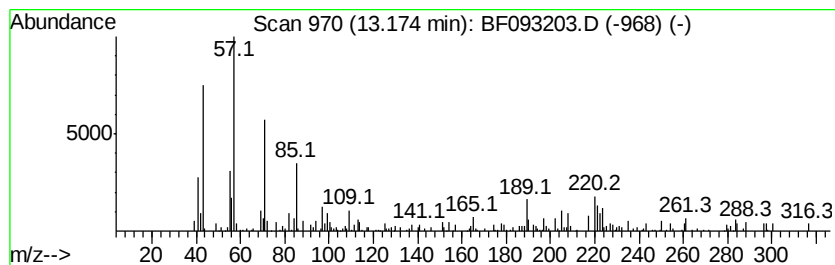
Instrument :  
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ClientSampleId :  
CLP-B-01(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.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 Tetradecane Concentration Rank 7

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.           |
|---------|---------|-------------------------------------|------------------|----------------|
| 13.17   | 2.62 ng | 151471                              | Chrysene-d12     | 14.01          |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |         | Tetradecane                         | 198 C14H30       | 000629-59-4 64 |
| 2       |         | Heptadecane, 2,6,10,15-tetramethyl- | 296 C21H44       | 054833-48-6 55 |
| 3       |         | Heneicosane                         | 296 C21H44       | 000629-94-7 55 |
| 4       |         | Pentadecane                         | 212 C15H32       | 000629-62-9 50 |
| 5       |         | Octadecane                          | 254 C18H38       | 000593-45-3 50 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\  
Data File : BF093203.D  
Acq On : 27 Feb 2017 12:37  
Operator : SJ/MA  
Sample : I1954-09  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

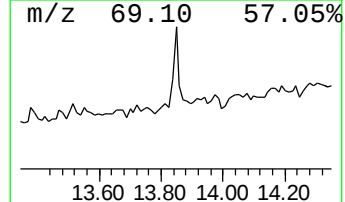
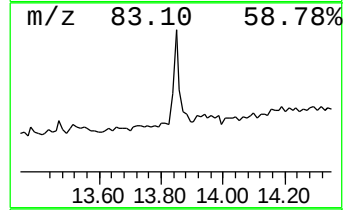
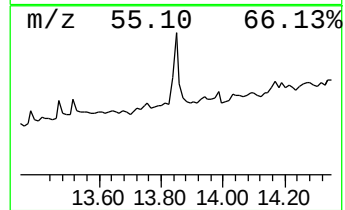
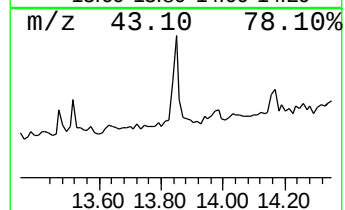
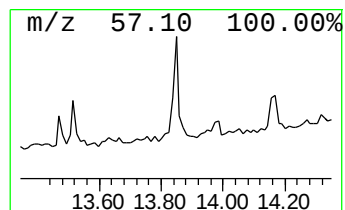
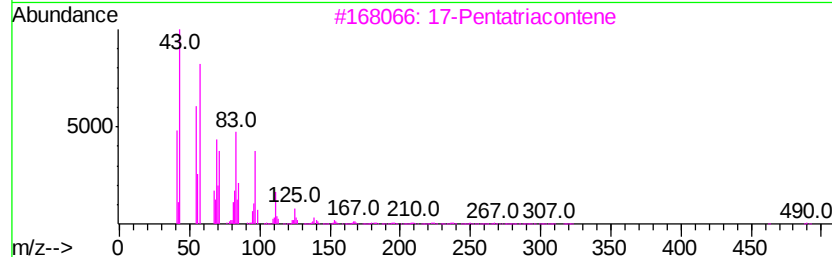
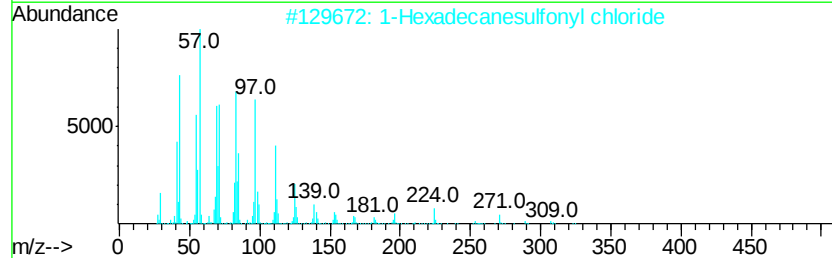
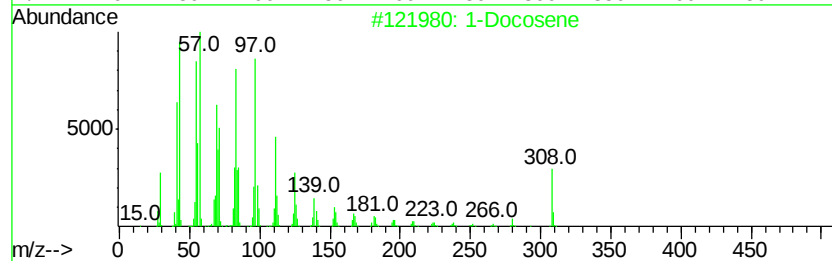
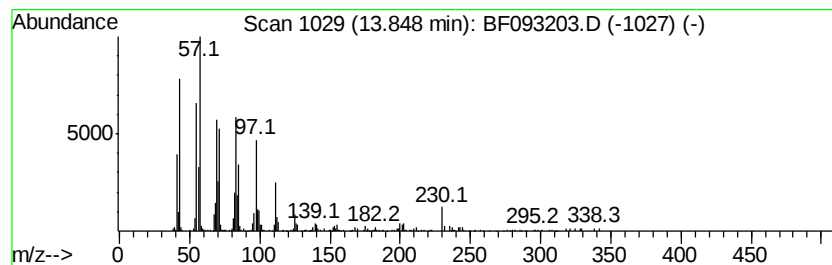
Instrument :  
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ClientSampleId :  
CLP-B-01(0-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.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-Docosene Concentration Rank 4

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.            |
|---------|-------------------------------------|--------------|------------------|-----------------|
| 13.85   | 6.39 ng                             | 369963       | Chrysene-d12     | 14.01           |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual       |
| 1       | 1-Docosene                          |              | 308 C22H44       | 001599-67-3 97  |
| 2       | 1-Hexadecanesulfonyl chloride       |              | 324 C16H33ClO2S  | 038775-38-1 91  |
| 3       | 17-Pentatriacontene                 |              | 491 C35H70       | 006971-40-0 87  |
| 4       | Pentafluoropropionic acid, hepta... |              | 402 C20H35F5O2   | 1000283-04-2 87 |
| 5       | Octadecane, 1-(ethenyloxy)-         |              | 296 C20H40O      | 000930-02-9 86  |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF022717\  
Data File : BF093203.D  
Acq On : 27 Feb 2017 12:37  
Operator : SJ/MA  
Sample : I1954-09  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CLP-B-01(0-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF013017.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... | 4.99  | 37.8    | ng    | 1930380  | 1                     | 6.83  | 1020730 | 20.0 |
| unknown6.60          | 6.60  | 86.6    | ng    | 4420350  | 1                     | 6.83  | 1020730 | 20.0 |
| Hexadecane           | 10.32 | 2.3     | ng    | 154691   | 3                     | 9.88  | 1373040 | 20.0 |
| Tridecane            | 10.78 | 4.9     | ng    | 303195   | 4                     | 11.37 | 1242170 | 20.0 |
| Pentadecane          | 11.23 | 2.3     | ng    | 139785   | 4                     | 11.37 | 1242170 | 20.0 |
| 11H-Benzo[b]fluorene | 13.13 | 2.7     | ng    | 155693   | 5                     | 14.01 | 1157770 | 20.0 |
| Tetradecane          | 13.17 | 2.6     | ng    | 151471   | 5                     | 14.01 | 1157770 | 20.0 |
| 1-Docosene           | 13.85 | 6.4     | ng    | 369963   | 5                     | 14.01 | 1157770 | 20.0 |