

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020116\  
 Data File : BF084731.D  
 Acq On : 2 Feb 2016 1:27  
 Operator : SJ/IZ  
 Sample : H1294-01  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EP002

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-BF020116.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         | 4.167       | 179           | 182         | 187          | rBV      | 226106         | 309497        | 6.02%           | 0.681%        |
| 2         | 4.876       | 240           | 244         | 246          | rBV      | 2474230        | 3928303       | 76.37%          | 8.640%        |
| 3         | 5.310       | 279           | 282         | 284          | rBV      | 4397472        | 4429141       | 86.11%          | 9.742%        |
| 4         | 5.710       | 315           | 317         | 319          | rBV      | 277929         | 237409        | 4.62%           | 0.522%        |
| 5         | 6.361       | 371           | 374         | 376          | rBV      | 4397771        | 4598146       | 89.39%          | 10.114%       |
| 6         | 6.476       | 382           | 384         | 387          | rBV      | 3773133        | 4113381       | 79.97%          | 9.048%        |
| 7         | 6.704       | 402           | 404         | 407          | rBV      | 882085         | 974053        | 18.94%          | 2.142%        |
| 8         | 6.864       | 416           | 418         | 421          | rVB      | 3679666        | 3202440       | 62.26%          | 7.044%        |
| 9         | 7.276       | 451           | 454         | 457          | rBV      | 2571443        | 2615710       | 50.85%          | 5.753%        |
| 10        | 7.996       | 514           | 517         | 519          | rBV      | 1550098        | 1339081       | 26.03%          | 2.945%        |
| 11        | 9.082       | 609           | 612         | 614          | rBV      | 4126640        | 5143660       | 100.00%         | 11.314%       |
| 12        | 9.470       | 644           | 646         | 648          | rBV      | 523494         | 480855        | 9.35%           | 1.058%        |
| 13        | 9.688       | 663           | 665         | 668          | rBV      | 73443          | 77362         | 1.50%           | 0.170%        |
| 14        | 9.745       | 668           | 670         | 673          | rVB      | 1553135        | 1440421       | 28.00%          | 3.168%        |
| 15        | 10.190      | 707           | 709         | 711          | rVB      | 127558         | 100333        | 1.95%           | 0.221%        |
| 16        | 10.408      | 725           | 728         | 731          | rBV      | 38048          | 59431         | 1.16%           | 0.131%        |
| 17        | 10.533      | 736           | 739         | 742          | rBV      | 3112527        | 3125479       | 60.76%          | 6.875%        |
| 18        | 10.659      | 748           | 750         | 755          | rVB      | 115303         | 131245        | 2.55%           | 0.289%        |
| 19        | 11.105      | 787           | 789         | 795          | rVB      | 50980          | 67627         | 1.31%           | 0.149%        |
| 20        | 11.231      | 797           | 800         | 802          | rBV      | 1183587        | 1448373       | 28.16%          | 3.186%        |
| 21        | 12.819      | 936           | 939         | 941          | rBV      | 3799149        | 4688829       | 91.16%          | 10.313%       |
| 22        | 13.745      | 1016          | 1020        | 1027         | rBV      | 53721          | 147356        | 2.86%           | 0.324%        |
| 23        | 13.859      | 1027          | 1030        | 1032         | rBV      | 1047940        | 1227883       | 23.87%          | 2.701%        |
| 24        | 14.705      | 1102          | 1104        | 1107         | rBV      | 60090          | 63573         | 1.24%           | 0.140%        |
| 25        | 15.231      | 1148          | 1150        | 1153         | rBV      | 569102         | 860595        | 16.73%          | 1.893%        |
| 26        | 15.277      | 1153          | 1154        | 1158         | rVB      | 50759          | 71819         | 1.40%           | 0.158%        |
| 27        | 15.665      | 1186          | 1188        | 1191         | rBV      | 65018          | 90075         | 1.75%           | 0.198%        |
| 28        | 16.111      | 1224          | 1227        | 1232         | rVB      | 69752          | 113796        | 2.21%           | 0.250%        |
| 29        | 16.625      | 1269          | 1272        | 1279         | rVB      | 52189          | 114773        | 2.23%           | 0.252%        |
| 30        | 17.243      | 1322          | 1326        | 1331         | rVB      | 46071          | 103926        | 2.02%           | 0.229%        |
| 31        | 17.963      | 1385          | 1389        | 1396         | rBV      | 29304          | 82721         | 1.61%           | 0.182%        |
| 32        | 18.820      | 1459          | 1464        | 1471         | rBV3     | 21683          | 76738         | 1.49%           | 0.169%        |

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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-BF020116.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

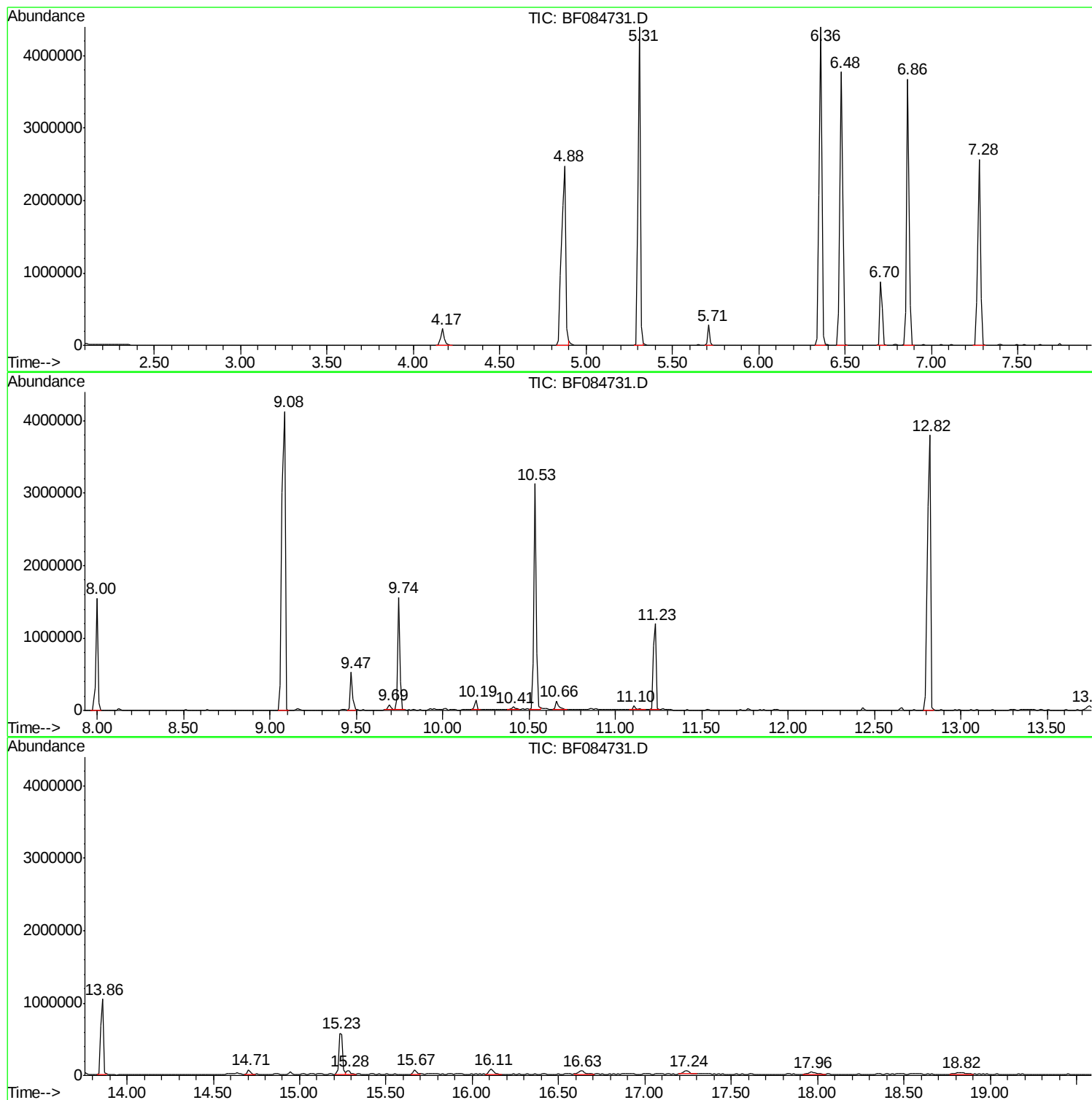
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
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TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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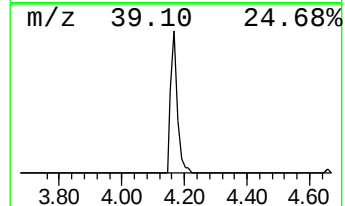
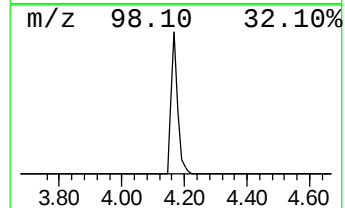
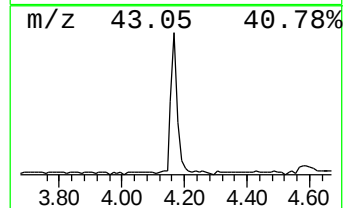
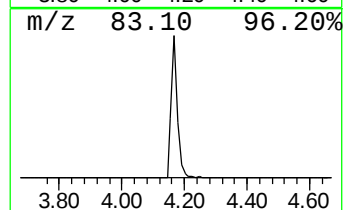
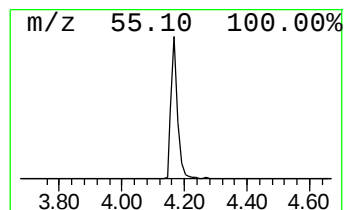
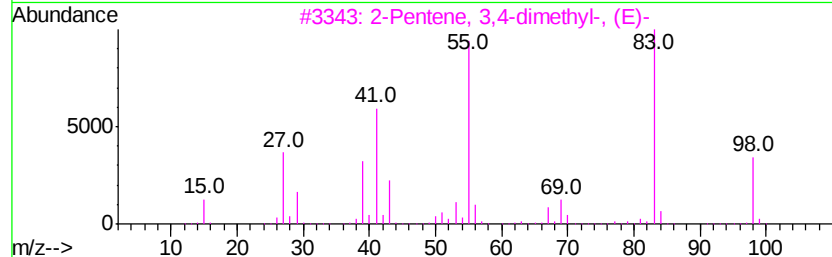
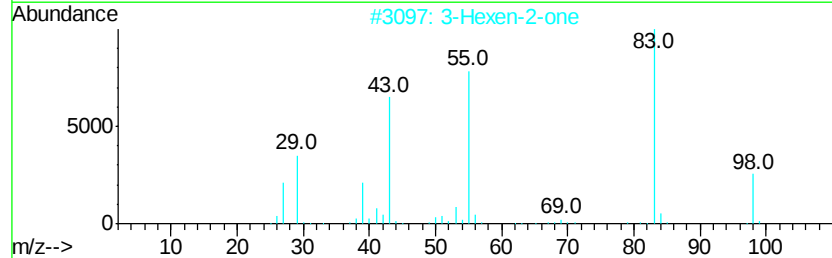
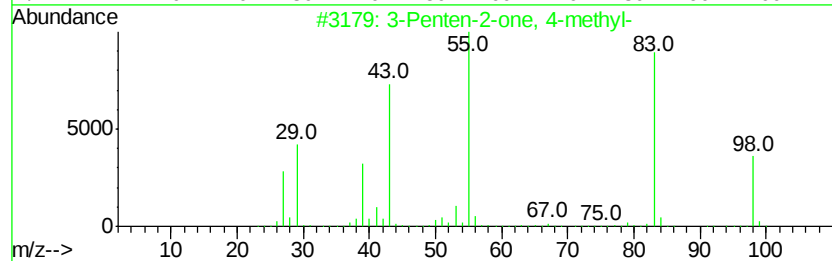
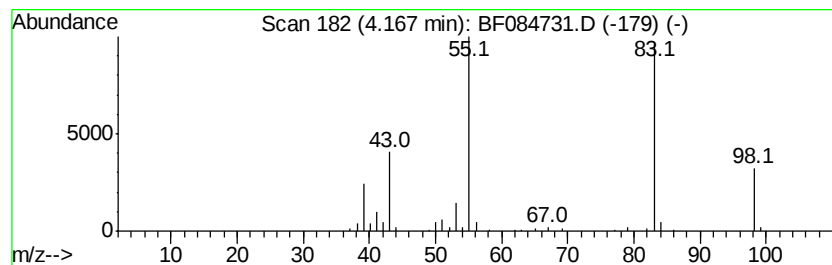
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 4

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.17 | 6.35 ng | 309497 | 1,4-Dichlorobenzene-d4 | 6.70 |

| Hit# | of | 5 | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|----|---------|-------------|------|
| 1    |    |   | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 91   |
| 2    |    |   | 3-Hexen-2-one                  | 98 | C6H10O  | 000763-93-9 | 90   |
| 3    |    |   | 2-Pentene, 3,4-dimethyl-, (E)- | 98 | C7H14   | 004914-92-5 | 86   |
| 4    |    |   | 2-Pentene, 2,4-dimethyl-       | 98 | C7H14   | 000625-65-0 | 86   |
| 5    |    |   | 2-Pentene, 3,4-dimethyl-, (Z)- | 98 | C7H14   | 004914-91-4 | 86   |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

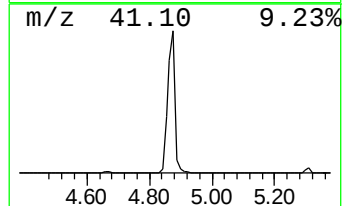
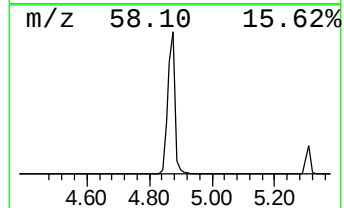
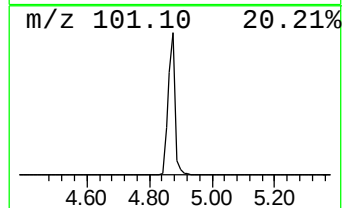
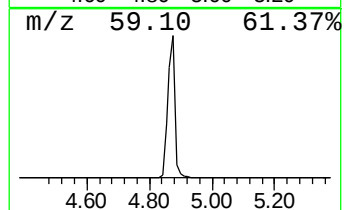
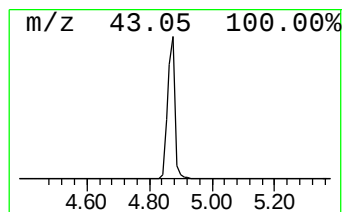
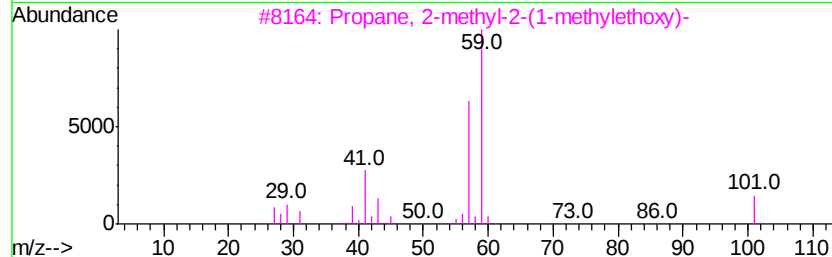
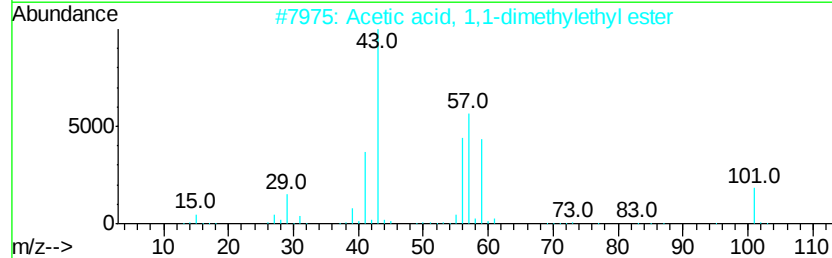
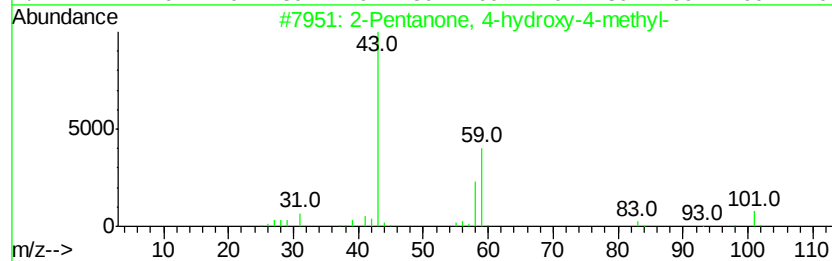
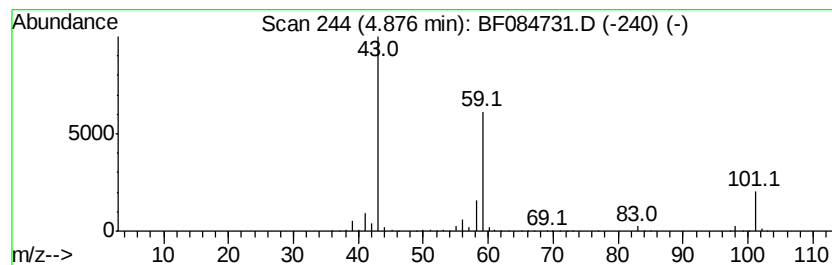
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 4.88 | 80.66 ng | 3928300 | 1,4-Dichlorobenzene-d4 | 6.70 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 39   |
| 3    |    | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O   | 017348-59-3 | 28   |
| 4    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 5    |    | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |



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Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP002

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

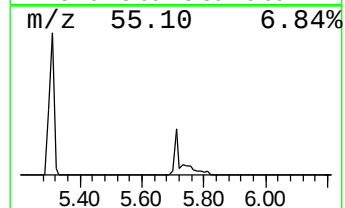
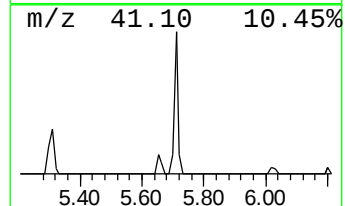
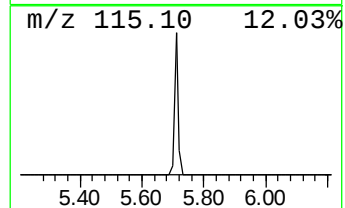
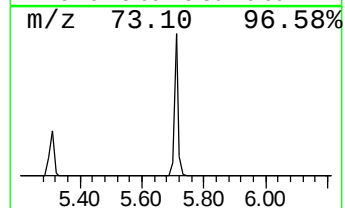
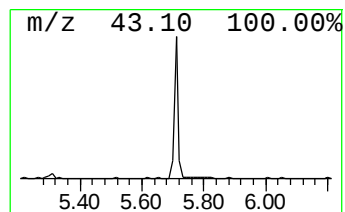
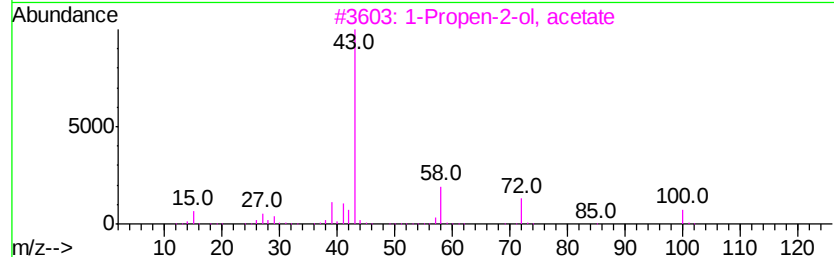
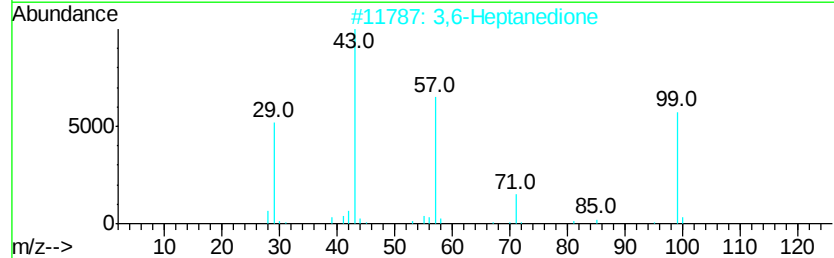
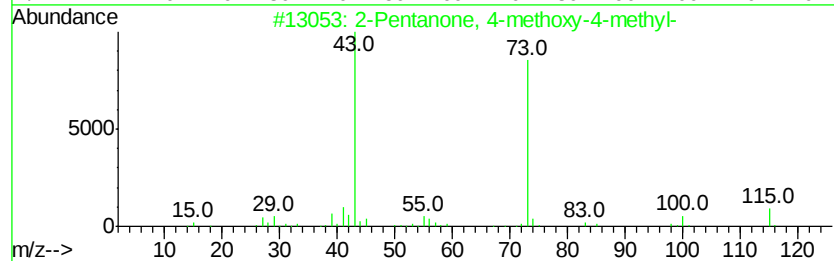
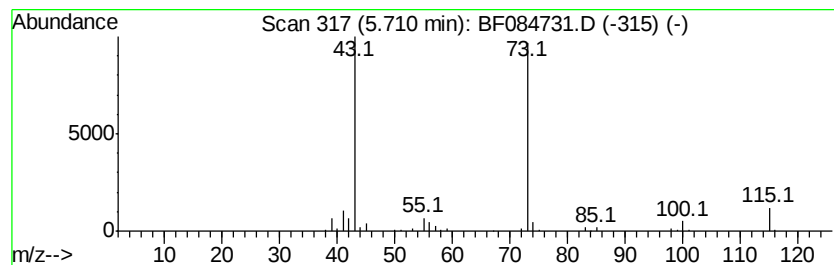
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 2-Pentanone, 4-methoxy-4-me... Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.71 | 4.87 ng | 237409 | 1,4-Dichlorobenzene-d4 | 6.70 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-methoxy-4-methyl-  | 130 | C7H14O2 | 000107-70-0 | 83   |
| 2    |    | 3,6-Heptanedione                  | 128 | C7H12O2 | 001703-51-1 | 17   |
| 3    |    | 1-Propen-2-ol, acetate            | 100 | C5H8O2  | 000108-22-5 | 10   |
| 4    |    | 1,3-Dioxolane, 2-(1-methylethyl)- | 116 | C6H12O2 | 000822-83-3 | 10   |
| 5    |    | Acetamide, N-butyl-               | 115 | C6H13NO | 001119-49-9 | 9    |



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ALS Vial : 29 Sample Multiplier: 1

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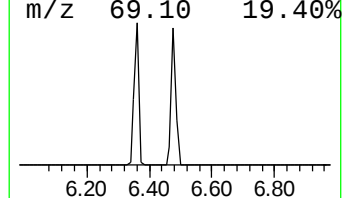
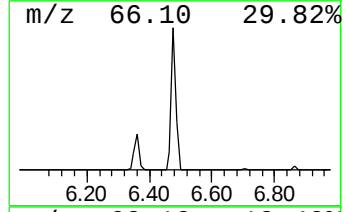
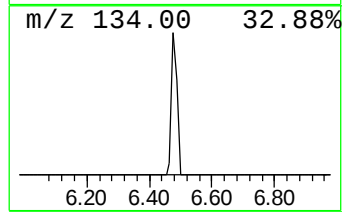
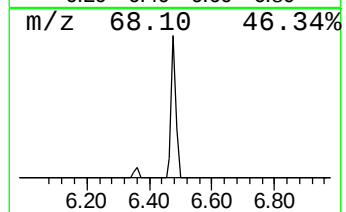
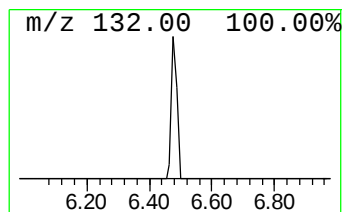
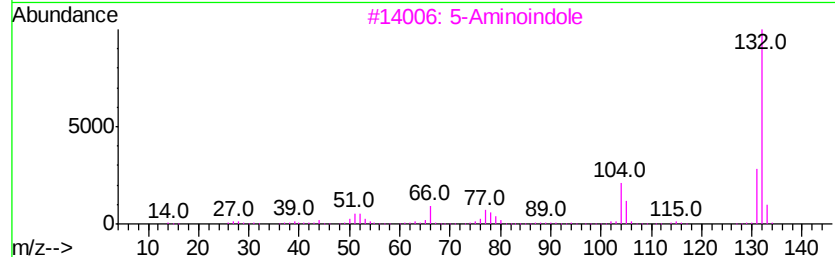
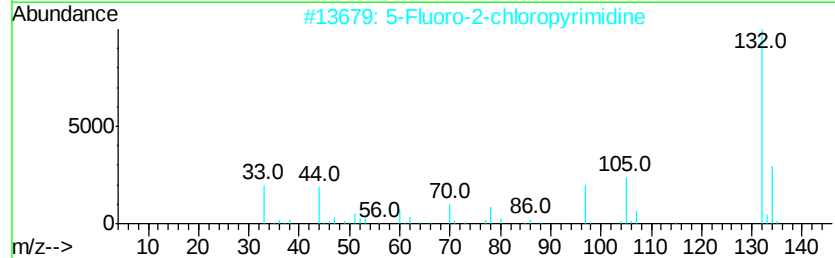
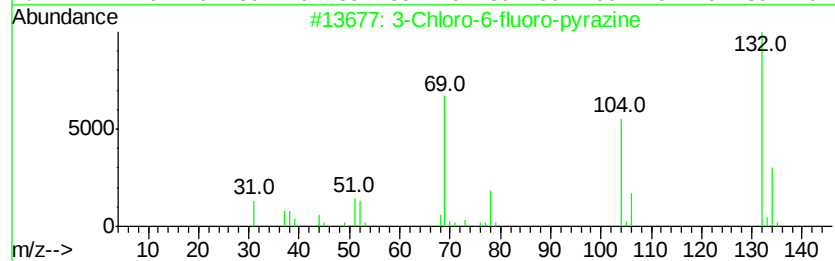
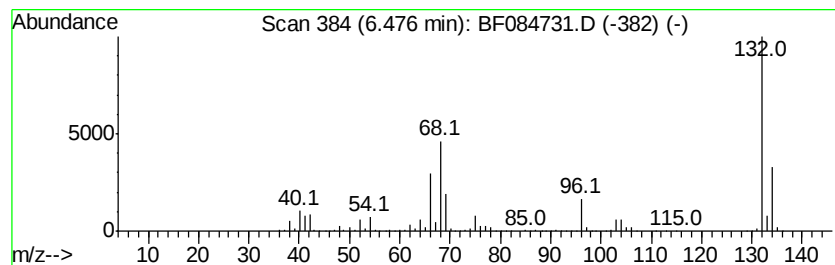
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 unknown6.48 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.48 | 84.46 ng | 4113380 | 1,4-Dichlorobenzene-d4 | 6.70 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0  | 12   |
| 3    |    | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | Tranylcypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |



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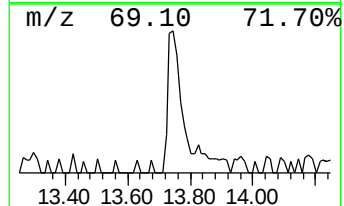
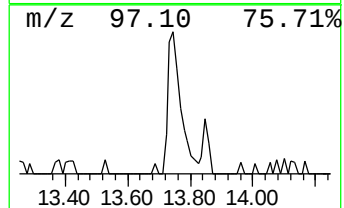
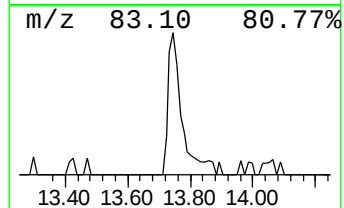
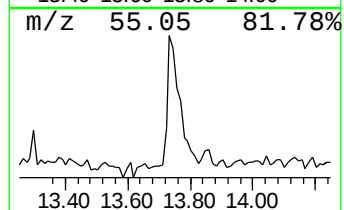
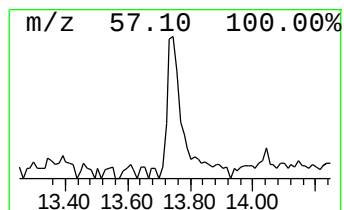
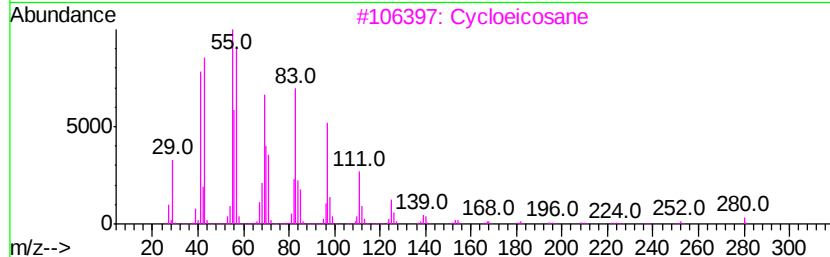
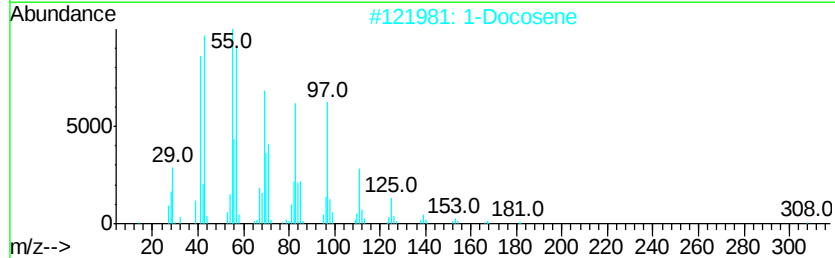
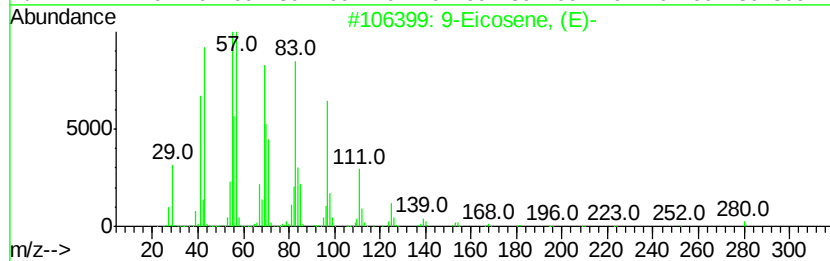
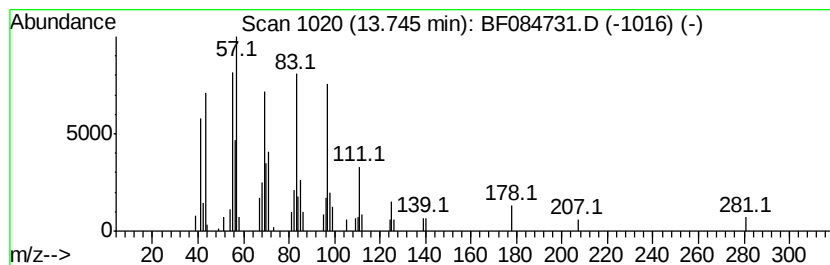
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 9-Eicosene, (E)- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.75 | 2.40 ng | 147356 | Chrysene-d12     | 13.86 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 9-Eicosene, (E)-                    | 280 | C20H40     | 074685-29-3  | 90   |
| 2    |    | 1-Docosene                          | 308 | C22H44     | 001599-67-3  | 74   |
| 3    |    | Cycloeicosane                       | 280 | C20H40     | 000296-56-0  | 68   |
| 4    |    | Heptafluorobutanoic acid, heptad... | 452 | C21H35F7O2 | 1000282-97-3 | 60   |
| 5    |    | Pentafluoropropionic acid, hepta... | 402 | C20H35F5O2 | 1000283-04-2 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020116\  
Data File : BF084731.D  
Acq On : 2 Feb 2016 1:27  
Operator : SJ/IZ  
Sample : H1294-01  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP002

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

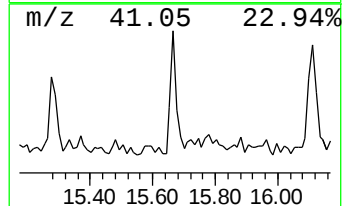
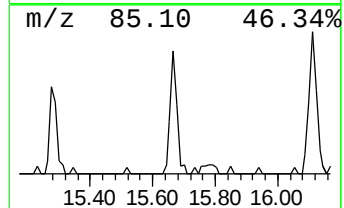
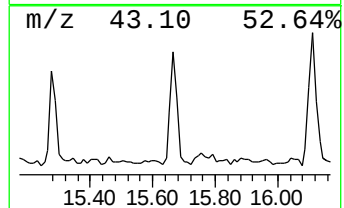
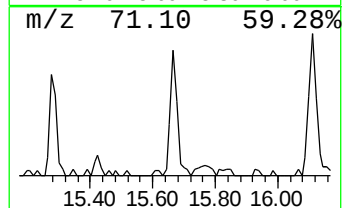
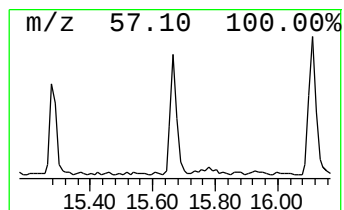
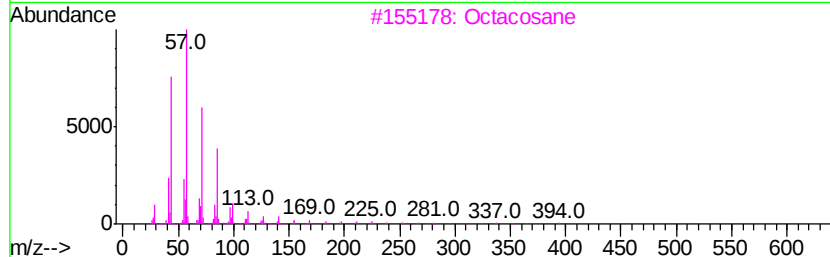
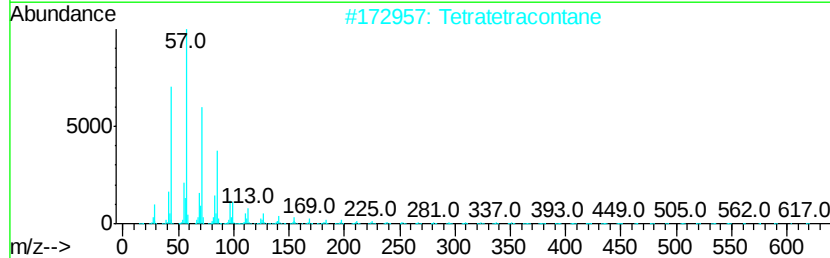
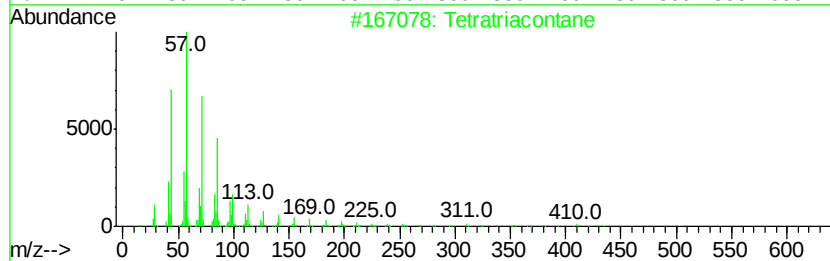
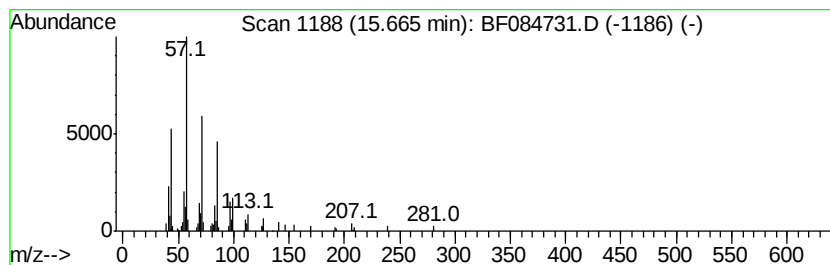
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Tetratriacontane Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.67 | 2.09 ng | 90075 | Perylene-d12     | 15.24 |

| Hit# of | 5 | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|---------|---|--------------------|-----|---------|-------------|------|
| 1       |   | Tetratriacontane   | 479 | C34H70  | 014167-59-0 | 91   |
| 2       |   | Tetratetracontane  | 619 | C44H90  | 007098-22-8 | 91   |
| 3       |   | Octacosane         | 394 | C28H58  | 000630-02-4 | 91   |
| 4       |   | Heneicosane        | 296 | C21H44  | 000629-94-7 | 90   |
| 5       |   | Eicosane, 9-octyl- | 394 | C28H58  | 013475-77-9 | 90   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020116\  
Data File : BF084731.D  
Acq On : 2 Feb 2016 1:27  
Operator : SJ/IZ  
Sample : H1294-01  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP002

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

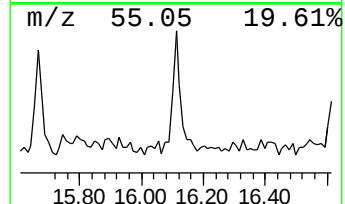
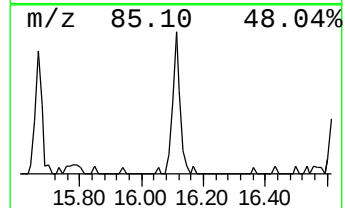
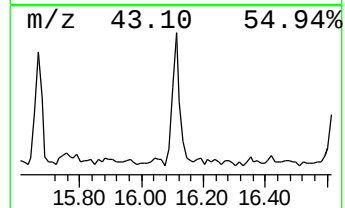
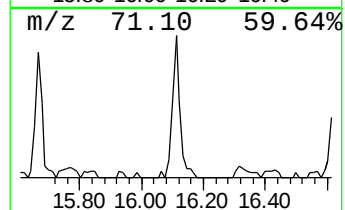
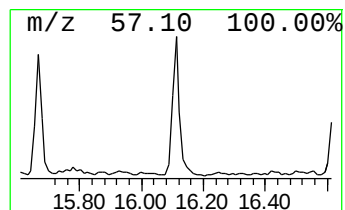
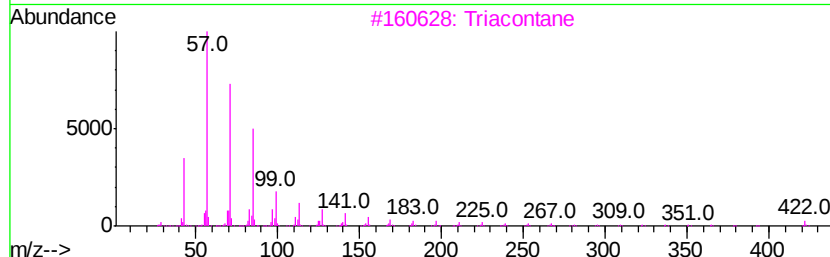
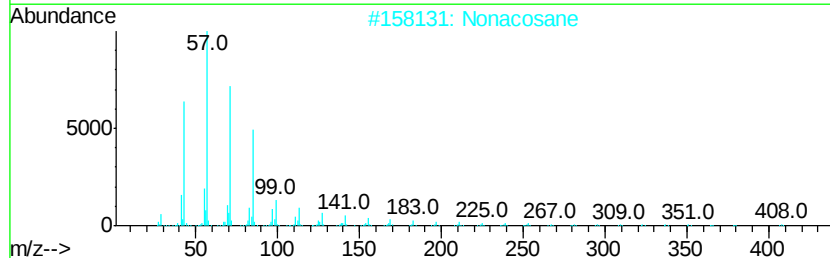
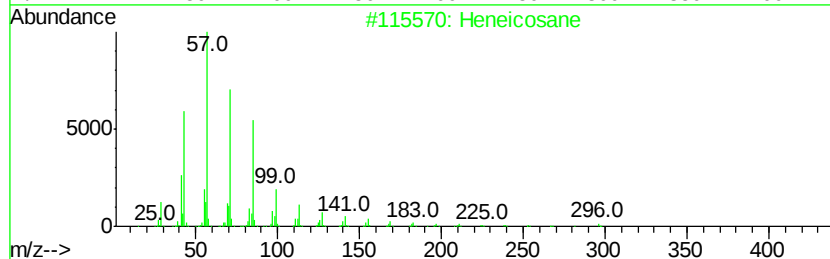
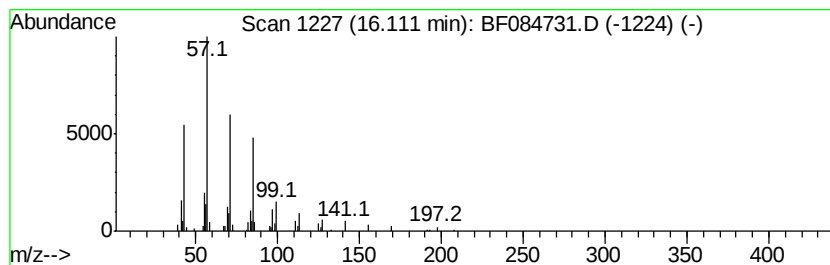
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Heneicosane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.11 | 2.64 ng | 113796 | Perylene-d12     | 15.24 |

| Hit# of | 5           | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------|--------------|-----|---------|-------------|------|
| 1       | Heneicosane |              | 296 | C21H44  | 000629-94-7 | 91   |
| 2       | Nonacosane  |              | 408 | C29H60  | 000630-03-5 | 91   |
| 3       | Triacontane |              | 422 | C30H62  | 000638-68-6 | 91   |
| 4       | Docosane    |              | 310 | C22H46  | 000629-97-0 | 86   |
| 5       | Octadecane  |              | 254 | C18H38  | 000593-45-3 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020116\  
Data File : BF084731.D  
Acq On : 2 Feb 2016 1:27  
Operator : SJ/IZ  
Sample : H1294-01  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP002

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

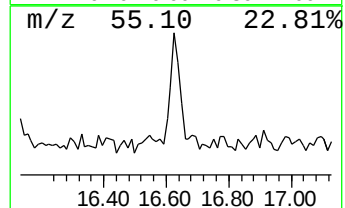
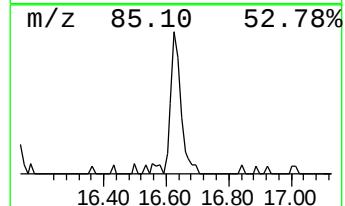
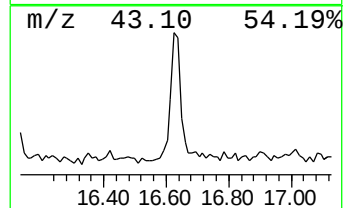
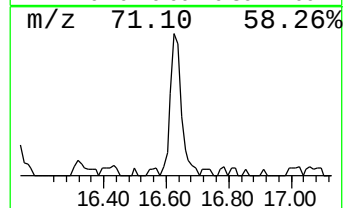
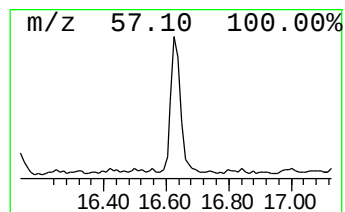
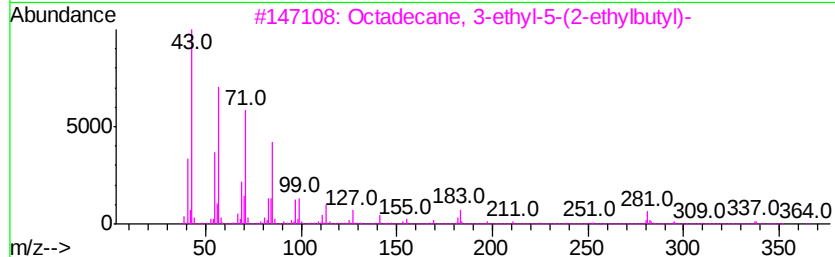
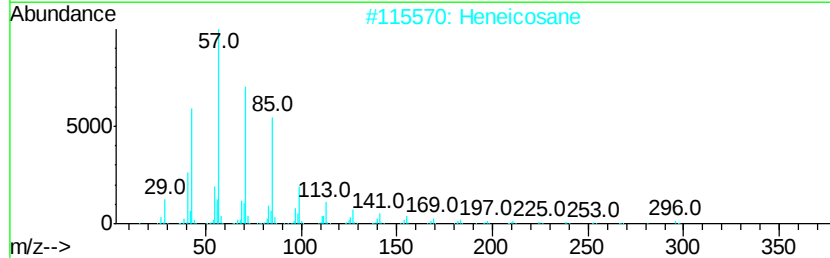
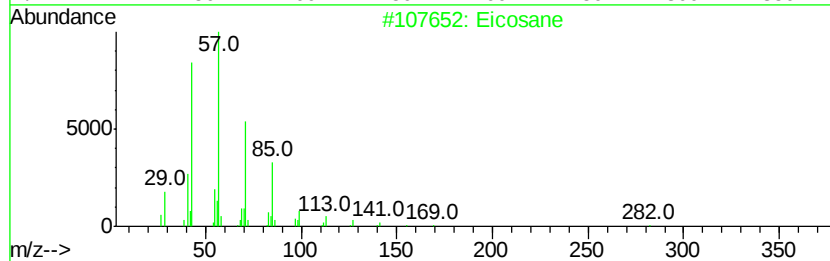
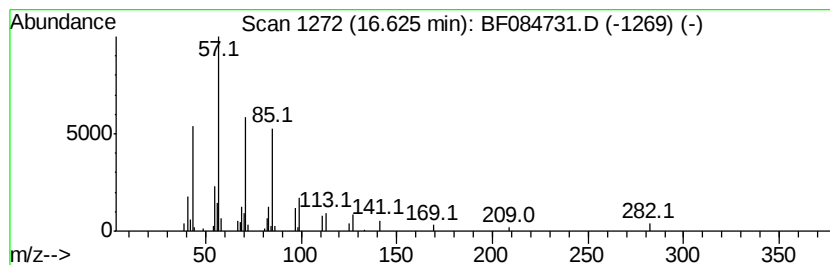
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Eicosane Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.63 | 2.67 ng | 114773 | Perylene-d12     | 15.24 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------------------------------|--------------|-----|---------|-------------|------|
| 1       | Eicosane                            |              | 282 | C20H42  | 000112-95-8 | 94   |
| 2       | Heneicosane                         |              | 296 | C21H44  | 000629-94-7 | 86   |
| 3       | Octadecane, 3-ethyl-5-(2-ethylbu... |              | 366 | C26H54  | 055282-12-7 | 78   |
| 4       | Hexatriacontane                     |              | 507 | C36H74  | 000630-06-8 | 78   |
| 5       | Tetracosane                         |              | 338 | C24H50  | 000646-31-1 | 72   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF020116\  
Data File : BF084731.D  
Acq On : 2 Feb 2016 1:27  
Operator : SJ/IZ  
Sample : H1294-01  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EP002

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF020116.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 3-Penten-2-one, 4... | 4.17  | 6.3     | ng    | 309497   | 1                     | 6.70  | 974053  | 20.0 |
| 2-Pentanone, 4-hy... | 4.88  | 80.7    | ng    | 3928300  | 1                     | 6.70  | 974053  | 20.0 |
| 2-Pentanone, 4-me... | 5.71  | 4.9     | ng    | 237409   | 1                     | 6.70  | 974053  | 20.0 |
| unknown6.48          | 6.48  | 84.5    | ng    | 4113380  | 1                     | 6.70  | 974053  | 20.0 |
| 9-Eicosene, (E)-     | 13.75 | 2.4     | ng    | 147356   | 5                     | 13.86 | 1227880 | 20.0 |
| Tetratriacontane     | 15.67 | 2.1     | ng    | 90075    | 6                     | 15.24 | 860595  | 20.0 |
| Heneicosane          | 16.11 | 2.6     | ng    | 113796   | 6                     | 15.24 | 860595  | 20.0 |
| Eicosane             | 16.63 | 2.7     | ng    | 114773   | 6                     | 15.24 | 860595  | 20.0 |