

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032216\  
 Data File : BF085695.D  
 Acq On : 23 Mar 2016 9:15  
 Operator : UM/SJ  
 Sample : H1857-09  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BLIND-DUPLICATE-1

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-BF031816.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         | 3.807       | 127           | 133         | 136          | rBV      | 4178260        | 10714052      | 100.00%         | 24.483%       |
| 2         | 4.493       | 190           | 193         | 196          | rBV      | 182660         | 212828        | 1.99%           | 0.486%        |
| 3         | 5.018       | 237           | 239         | 249          | rVB      | 267709         | 501427        | 4.68%           | 1.146%        |
| 4         | 5.441       | 273           | 276         | 278          | rBV      | 2980276        | 3523934       | 32.89%          | 8.053%        |
| 5         | 6.470       | 363           | 366         | 369          | rBV      | 2544488        | 3389591       | 31.64%          | 7.746%        |
| 6         | 6.607       | 375           | 378         | 380          | rBV      | 2988799        | 3795192       | 35.42%          | 8.673%        |
| 7         | 6.836       | 395           | 398         | 400          | rBV      | 563015         | 742501        | 6.93%           | 1.697%        |
| 8         | 6.984       | 409           | 411         | 414          | rBV      | 2359056        | 2674427       | 24.96%          | 6.111%        |
| 9         | 7.396       | 444           | 447         | 450          | rBV      | 1862194        | 2347049       | 21.91%          | 5.363%        |
| 10        | 8.116       | 507           | 510         | 512          | rBV      | 1256597        | 1069375       | 9.98%           | 2.444%        |
| 11        | 9.190       | 601           | 604         | 607          | rBV      | 3024868        | 3650596       | 34.07%          | 8.342%        |
| 12        | 9.590       | 637           | 639         | 641          | rBV      | 701689         | 729549        | 6.81%           | 1.667%        |
| 13        | 9.807       | 654           | 658         | 660          | rBV      | 129081         | 202567        | 1.89%           | 0.463%        |
| 14        | 9.865       | 661           | 663         | 665          | rVB      | 1008933        | 1106567       | 10.33%          | 2.529%        |
| 15        | 10.299      | 698           | 701         | 703          | rBV      | 209659         | 265947        | 2.48%           | 0.608%        |
| 16        | 10.425      | 709           | 712         | 714          | rBV2     | 53216          | 110170        | 1.03%           | 0.252%        |
| 17        | 10.516      | 717           | 720         | 724          | rBV2     | 94849          | 229698        | 2.14%           | 0.525%        |
| 18        | 10.665      | 729           | 733         | 735          | rBV2     | 1773325        | 2527113       | 23.59%          | 5.775%        |
| 19        | 10.779      | 741           | 743         | 747          | rVB      | 149578         | 300878        | 2.81%           | 0.688%        |
| 20        | 11.351      | 791           | 793         | 795          | rBV      | 1191914        | 1022608       | 9.54%           | 2.337%        |
| 21        | 11.876      | 838           | 839         | 845          | rBV5     | 91441          | 208667        | 1.95%           | 0.477%        |
| 22        | 12.665      | 906           | 908         | 912          | rVB2     | 84795          | 135707        | 1.27%           | 0.310%        |
| 23        | 12.939      | 929           | 932         | 934          | rBV      | 2462597        | 2826573       | 26.38%          | 6.459%        |
| 24        | 13.831      | 1008          | 1010        | 1018         | rVB      | 93145          | 166802        | 1.56%           | 0.381%        |
| 25        | 13.979      | 1021          | 1023        | 1026         | rBV      | 718088         | 654864        | 6.11%           | 1.496%        |
| 26        | 15.397      | 1144          | 1147        | 1153         | rVB      | 533636         | 652411        | 6.09%           | 1.491%        |

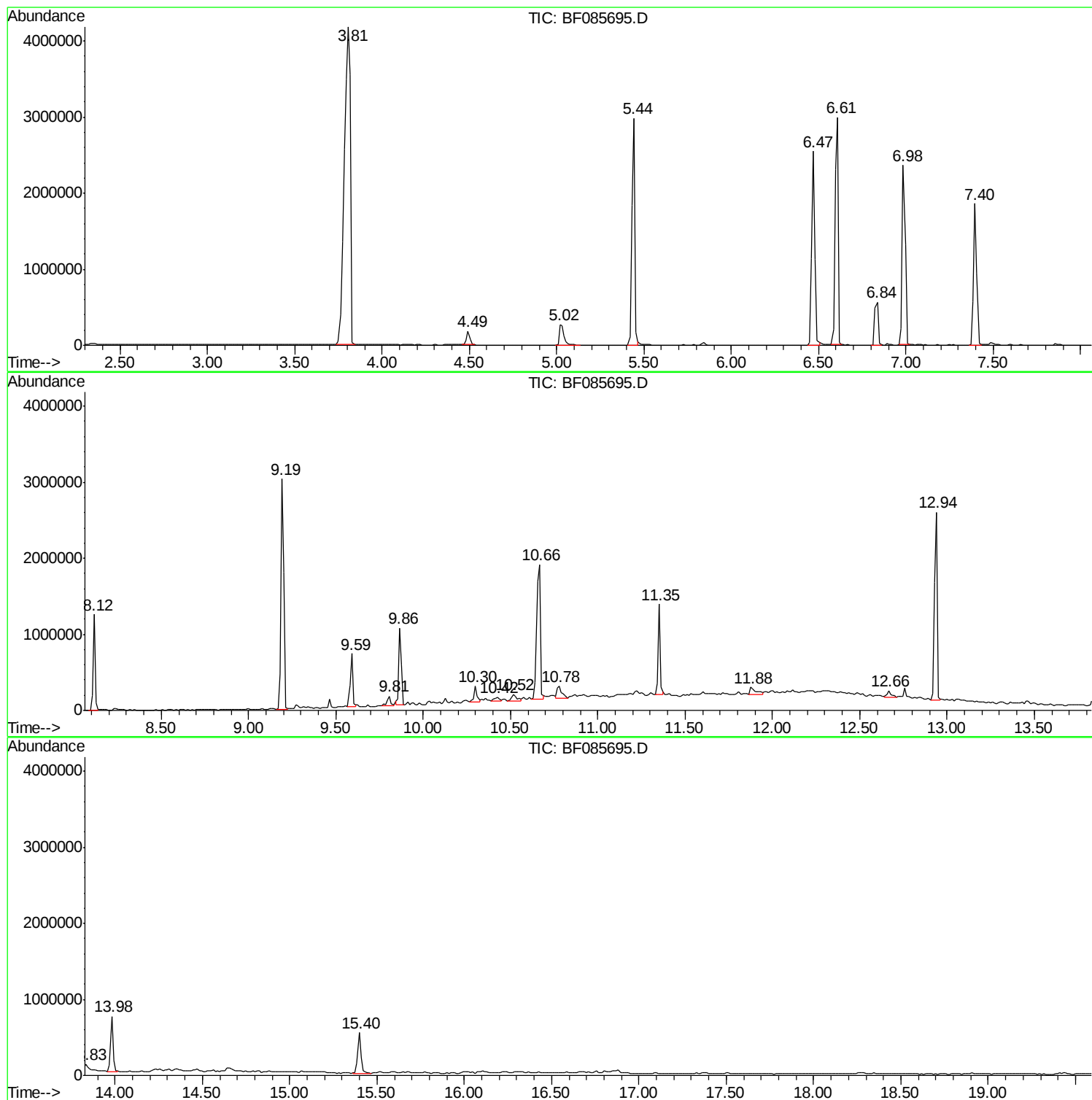
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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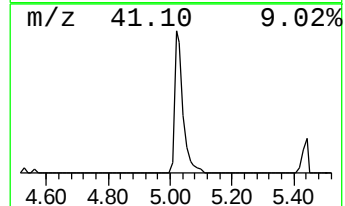
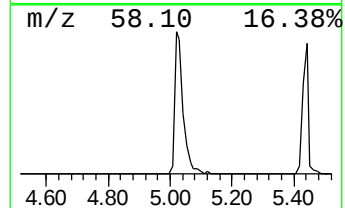
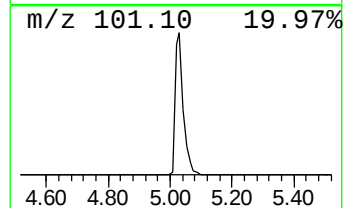
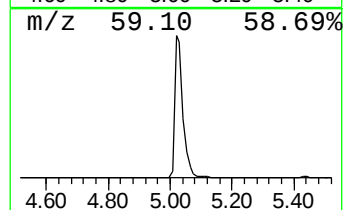
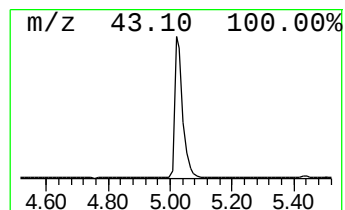
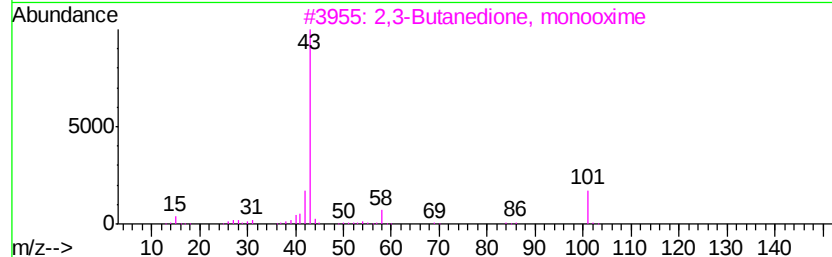
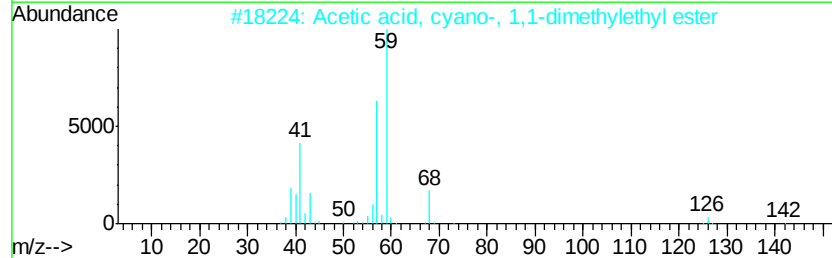
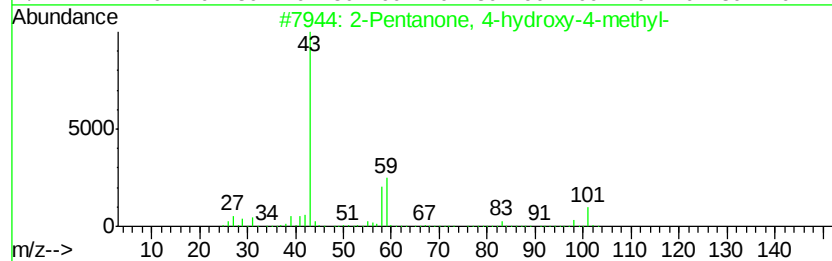
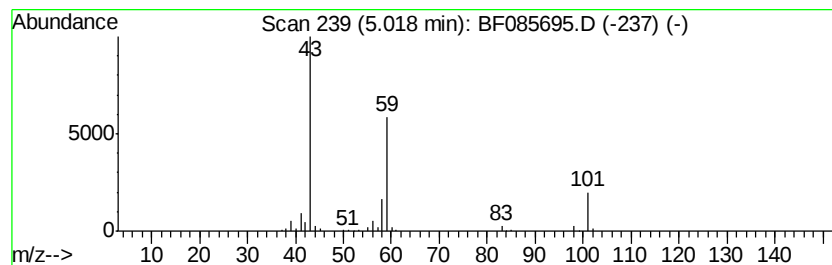
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.02 | 13.51 ng | 501427 | 1,4-Dichlorobenzene-d4 | 6.84 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|---------|---|-------------------------------------|-----|----------|-------------|------|
| 1       |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 64   |
| 2       |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 32   |
| 3       |   | 2,3-Butanedione, monooxime          | 101 | C4H7NO2  | 000057-71-6 | 25   |
| 4       |   | Butane, 1-ethoxy-                   | 102 | C6H14O   | 000628-81-9 | 9    |
| 5       |   | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

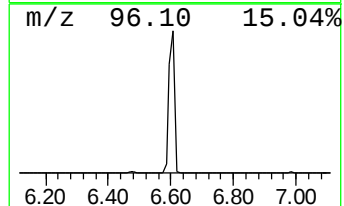
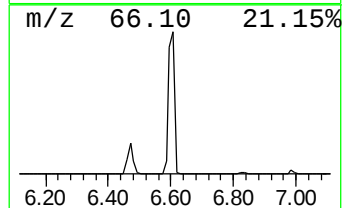
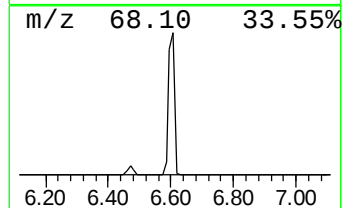
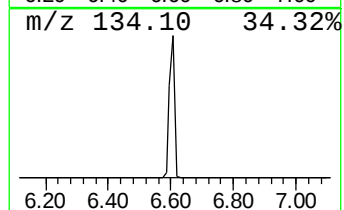
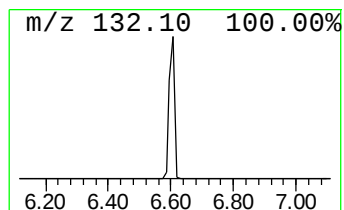
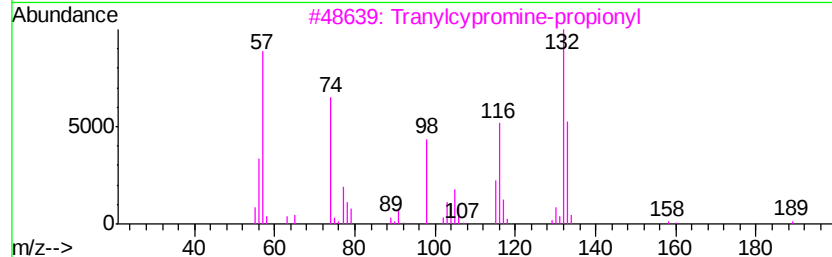
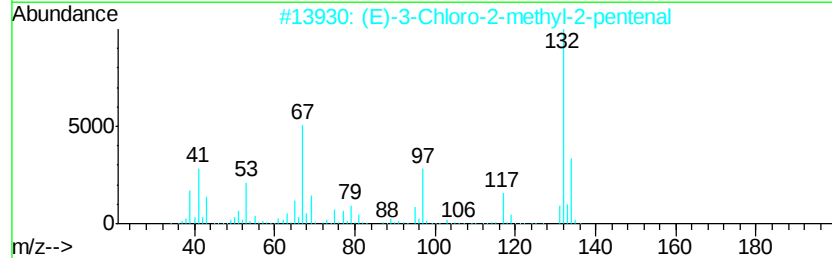
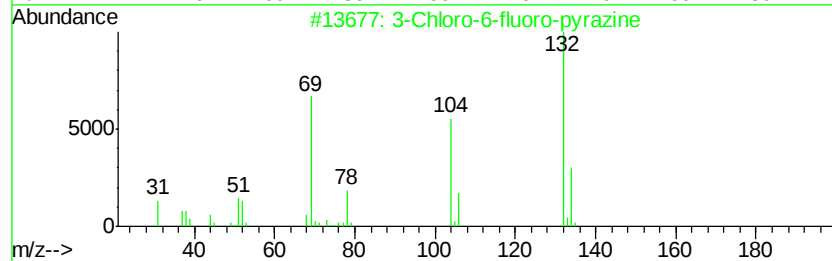
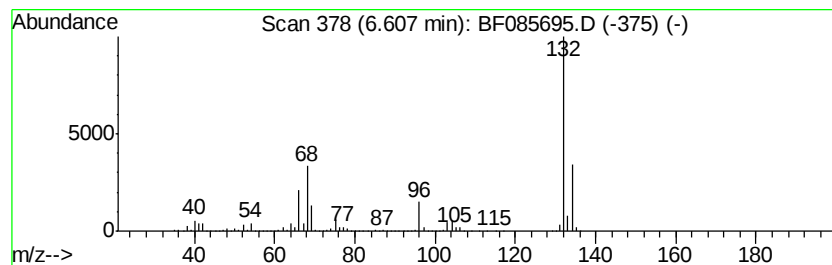
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown6.61 Concentration Rank 2

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 6.61 | 102.23 ng | 3795190 | 1,4-Dichlorobenzene-d4 | 6.84 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 16   |
| 4    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 12   |
| 5    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 9    |



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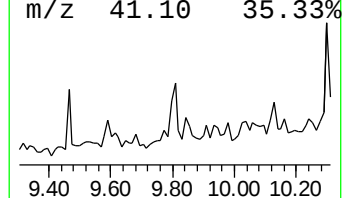
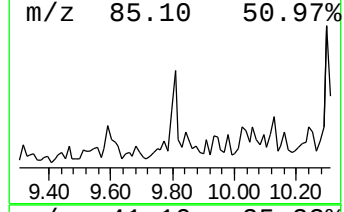
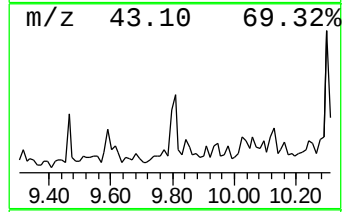
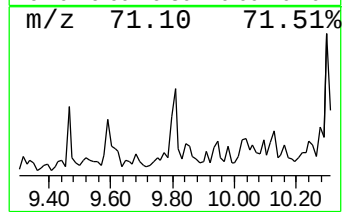
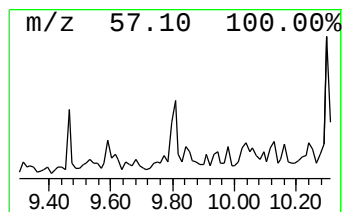
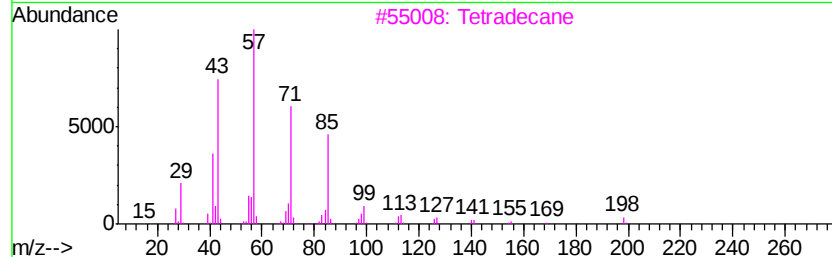
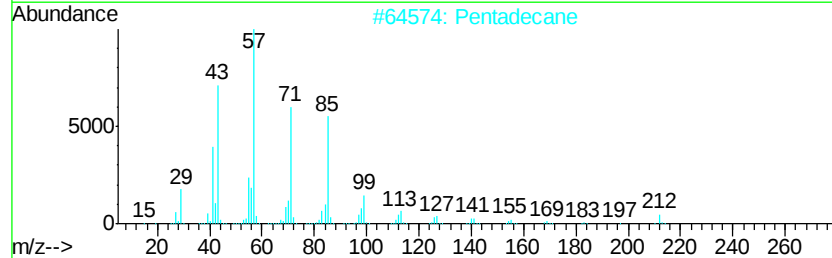
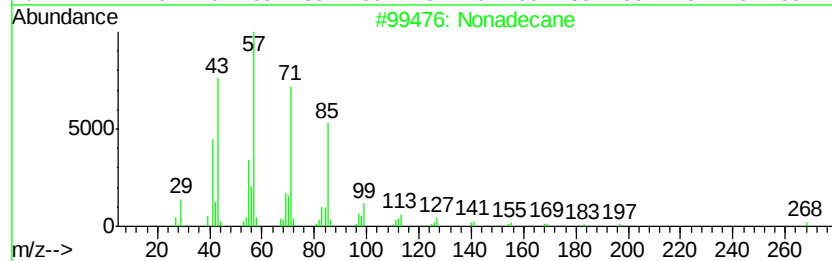
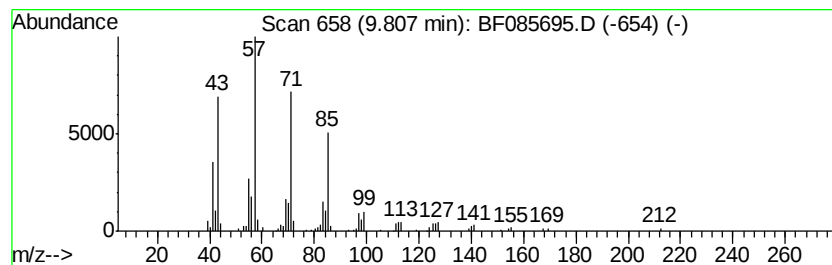
Instrument :  
BNA\_F  
ClientSampleId :  
BLIND-DUPLICATE-1

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

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Peak Number 6 Nonadecane Concentration Rank 11

| R.T.      | EstConc                | Area   | Relative to ISTD |             | R.T. |
|-----------|------------------------|--------|------------------|-------------|------|
| 9.81      | 3.66 ng                | 202567 | Acenaphthene-d10 |             | 9.86 |
| Hit# of 5 | Tentative ID           | MW     | MolForm          | CAS#        | Qual |
| 1         | Nonadecane             | 268    | C19H40           | 000629-92-5 | 90   |
| 2         | Pentadecane            | 212    | C15H32           | 000629-62-9 | 86   |
| 3         | Tetradecane            | 198    | C14H30           | 000629-59-4 | 83   |
| 4         | Heptadecane, 8-methyl- | 254    | C18H38           | 013287-23-5 | 80   |
| 5         | Tridecane              | 184    | C13H28           | 000629-50-5 | 80   |



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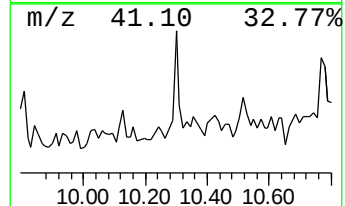
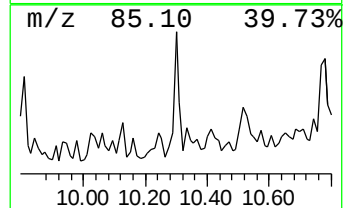
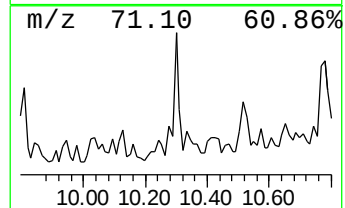
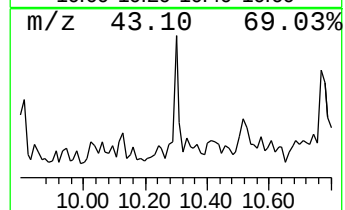
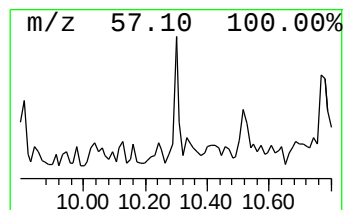
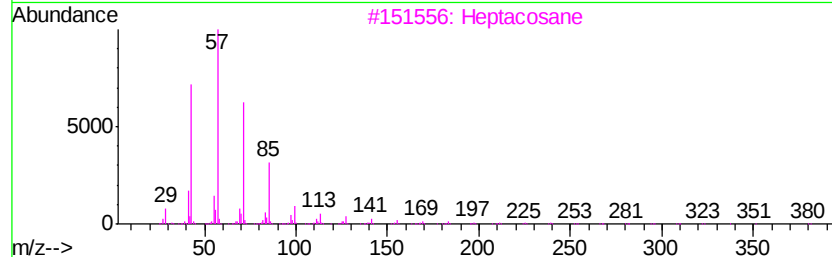
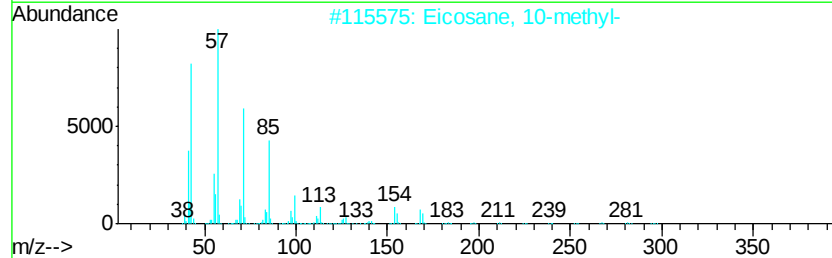
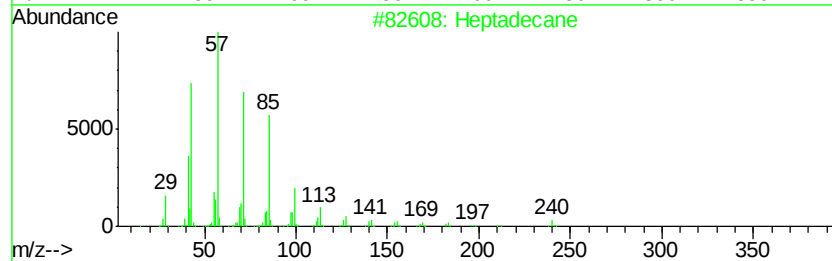
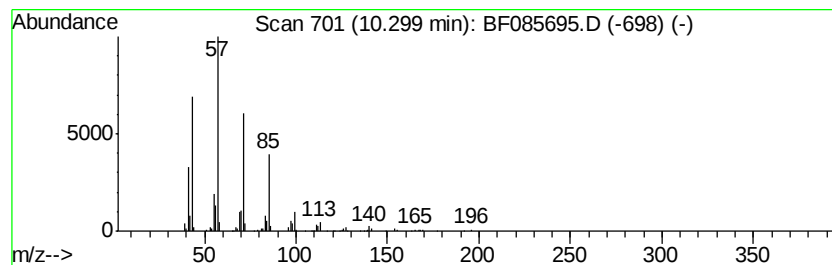
Instrument :  
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ClientSampleId :  
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Peak Number 7 Heptadecane Concentration Rank 8

| R.T.    | EstConc | Area                 | Relative to ISTD |         | R.T.        |      |
|---------|---------|----------------------|------------------|---------|-------------|------|
| 10.30   | 4.81 ng | 265947               | Acenaphthene-d10 |         | 9.86        |      |
| Hit# of | 5       | Tentative ID         | MW               | MolForm | CAS#        | Qual |
| 1       |         | Heptadecane          | 240              | C17H36  | 000629-78-7 | 90   |
| 2       |         | Eicosane, 10-methyl- | 296              | C21H44  | 054833-23-7 | 90   |
| 3       |         | Heptacosane          | 380              | C27H56  | 000593-49-7 | 90   |
| 4       |         | Nonadecane           | 268              | C19H40  | 000629-92-5 | 87   |
| 5       |         | Tetracosane          | 338              | C24H50  | 000646-31-1 | 86   |



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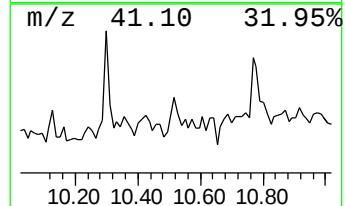
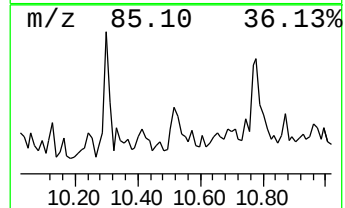
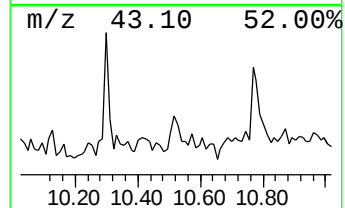
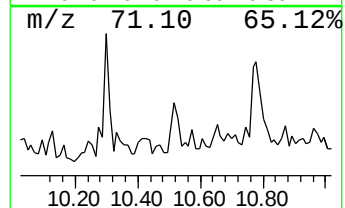
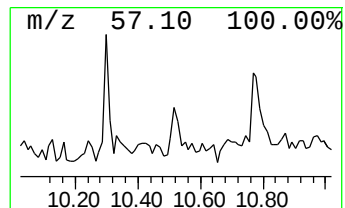
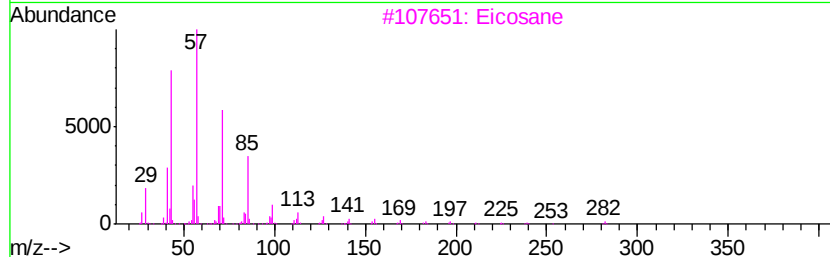
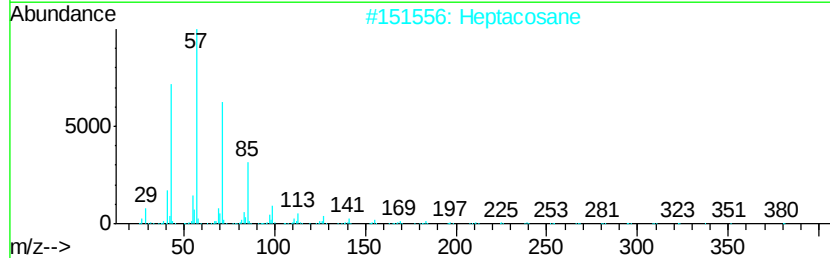
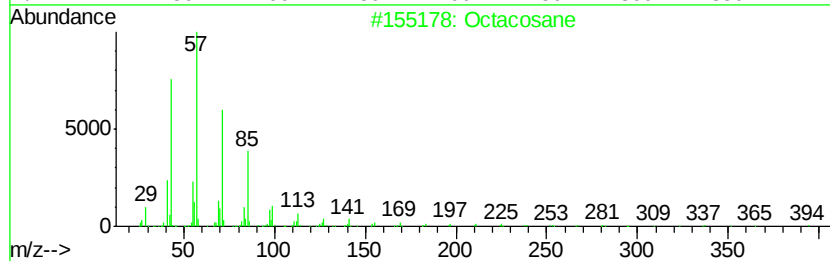
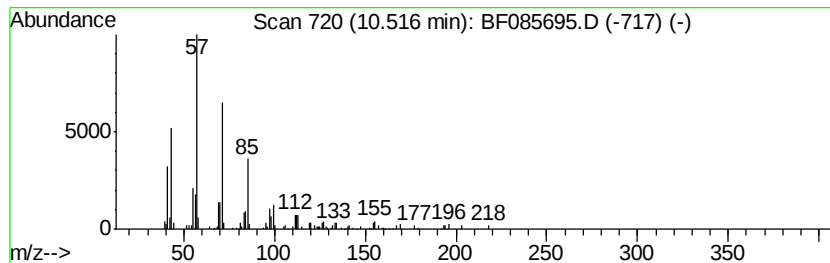
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Octacosane Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T. |
|-------|---------|--------|------------------|------|
| 10.52 | 4.15 ng | 229698 | Acenaphthene-d10 | 9.86 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Octacosane        | 394 | C28H58  | 000630-02-4 | 86   |
| 2    |    | Heptacosane       | 380 | C27H56  | 000593-49-7 | 86   |
| 3    |    | Eicosane          | 282 | C20H42  | 000112-95-8 | 86   |
| 4    |    | Hexadecane        | 226 | C16H34  | 000544-76-3 | 80   |
| 5    |    | Tetratetracontane | 619 | C44H90  | 007098-22-8 | 80   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032216\  
Data File : BF085695.D  
Acq On : 23 Mar 2016 9:15  
Operator : UM/SJ  
Sample : H1857-09  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

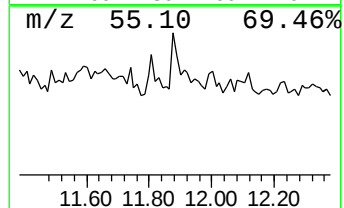
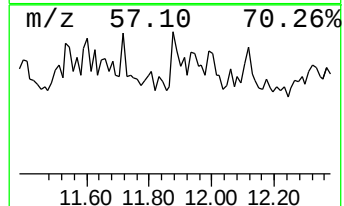
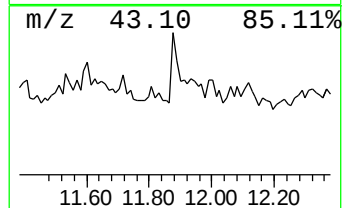
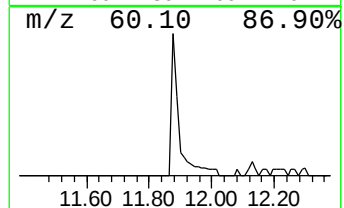
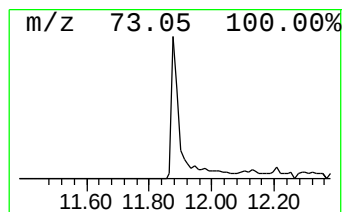
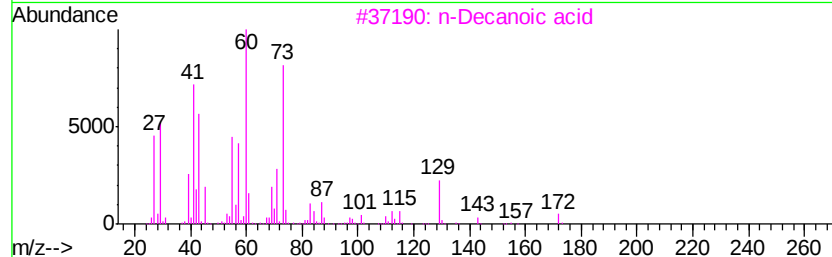
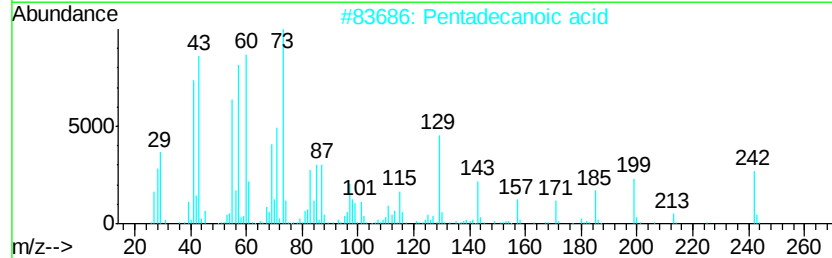
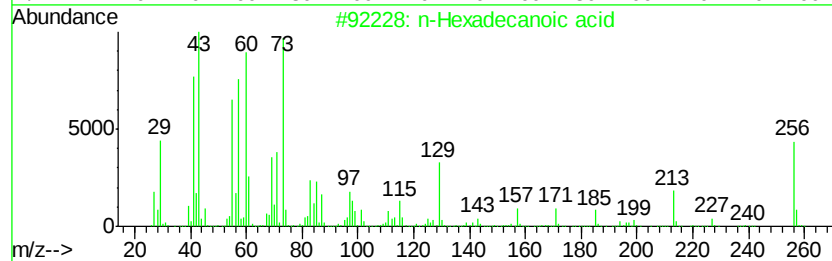
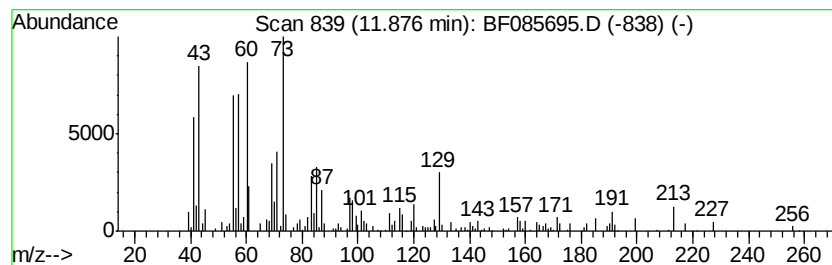
Instrument :  
BNA\_F  
ClientSampleId :  
BLIND-DUPLICATE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 n-Hexadecanoic acid Concentration Rank 10

| R.T.    | EstConc | Area                | Relative to ISTD |          | R.T.        |      |
|---------|---------|---------------------|------------------|----------|-------------|------|
| 11.88   | 4.08 ng | 208667              | Phenanthrene-d10 |          | 11.35       |      |
| Hit# of | 5       | Tentative ID        | MW               | MolForm  | CAS#        | Qual |
| 1       |         | n-Hexadecanoic acid | 256              | C16H32O2 | 000057-10-3 | 95   |
| 2       |         | Pentadecanoic acid  | 242              | C15H30O2 | 001002-84-2 | 90   |
| 3       |         | n-Decanoic acid     | 172              | C10H20O2 | 000334-48-5 | 68   |
| 4       |         | Nonanoic acid       | 158              | C9H18O2  | 000112-05-0 | 18   |
| 5       |         | Octanoic Acid       | 144              | C8H16O2  | 000124-07-2 | 18   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032216\  
Data File : BF085695.D  
Acq On : 23 Mar 2016 9:15  
Operator : UM/SJ  
Sample : H1857-09  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

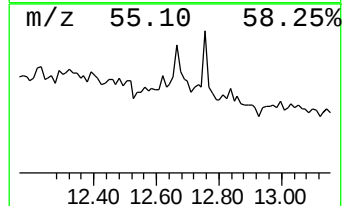
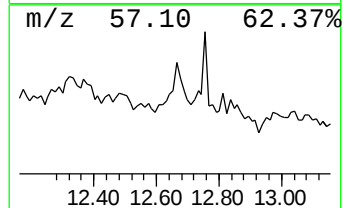
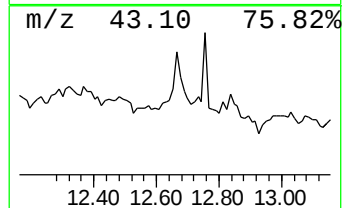
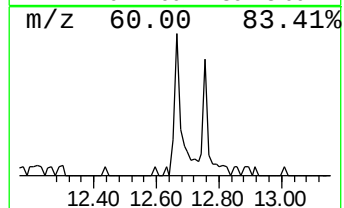
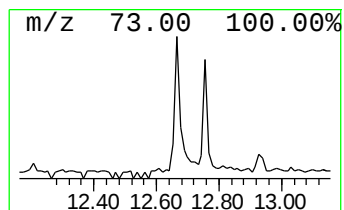
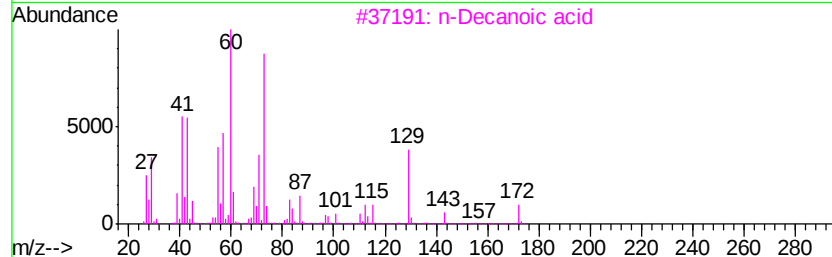
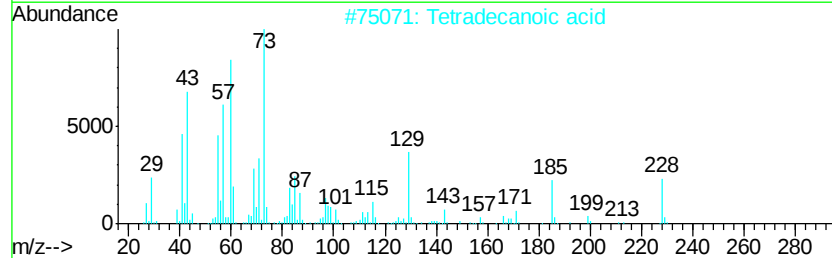
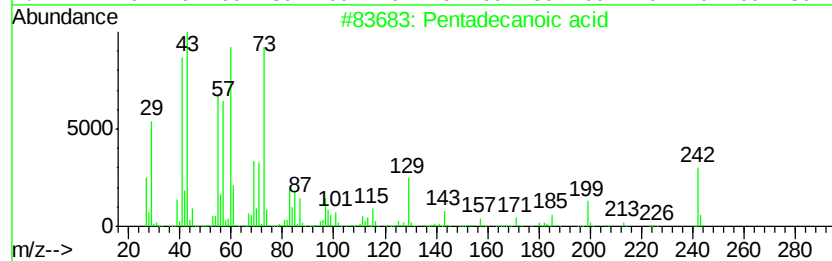
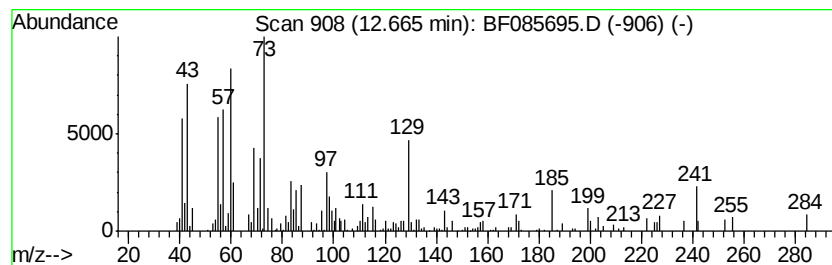
Instrument :  
BNA\_F  
ClientSampleId :  
BLIND-DUPLICATE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.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 11 Pentadecanoic acid Concentration Rank 12

| R.T.    | EstConc | Area               | Relative to ISTD |          | R.T.        |      |
|---------|---------|--------------------|------------------|----------|-------------|------|
| 12.66   | 2.65 ng | 135707             | Phenanthrene-d10 |          | 11.35       |      |
| Hit# of | 5       | Tentative ID       | MW               | MolForm  | CAS#        | Qual |
| 1       |         | Pentadecanoic acid | 242              | C15H30O2 | 001002-84-2 | 92   |
| 2       |         | Tetradecanoic acid | 228              | C14H28O2 | 000544-63-8 | 72   |
| 3       |         | n-Decanoic acid    | 172              | C10H20O2 | 000334-48-5 | 64   |
| 4       |         | Tridecanoic acid   | 214              | C13H26O2 | 000638-53-9 | 64   |
| 5       |         | Dodecanoic acid    | 200              | C12H24O2 | 000143-07-7 | 55   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF032216\  
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Operator : UM/SJ  
Sample : H1857-09  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BLIND-DUPLICATE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

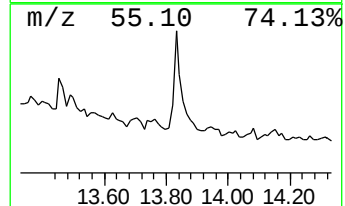
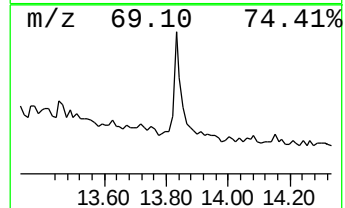
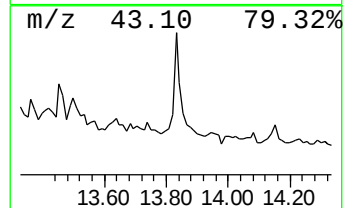
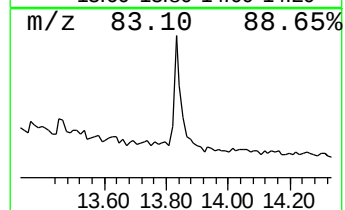
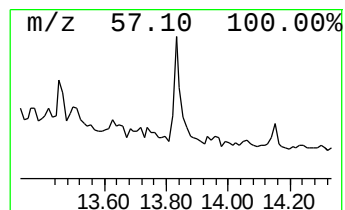
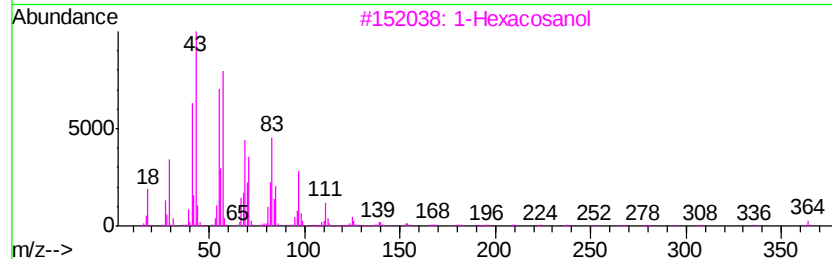
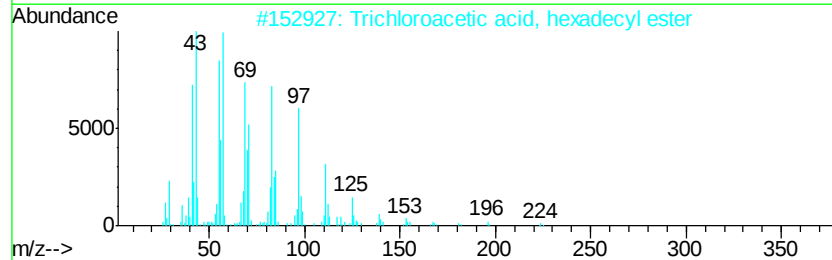
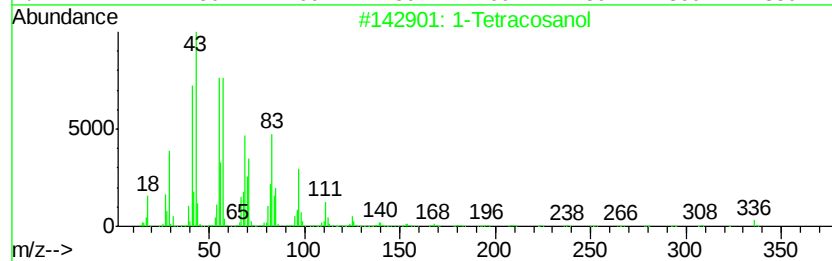
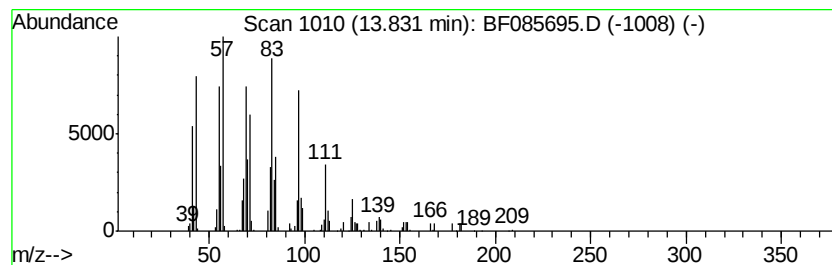
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 1-Tetracosanol Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.83 | 5.09 ng | 166802 | Chrysene-d12     | 13.98 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 1-Tetracosanol                      | 354 | C24H50O     | 000506-51-4  | 91   |
| 2    |    |   | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1  | 91   |
| 3    |    |   | 1-Hexacosanol                       | 382 | C26H54O     | 000506-52-5  | 91   |
| 4    |    |   | Pentafluoropropionic acid, octad... | 416 | C21H37F5O2  | 1000280-07-7 | 90   |
| 5    |    |   | 17-Pentatriacontene                 | 491 | C35H70      | 006971-40-0  | 90   |



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Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BLIND-DUPLICATE-1

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF031816.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... | 5.02  | 13.5    | ng    | 501427   | 1                     | 6.84  | 742501  | 20.0 |
| unknown6.61          | 6.61  | 102.2   | ng    | 3795190  | 1                     | 6.84  | 742501  | 20.0 |
| Nonadecane           | 9.81  | 3.7     | ng    | 202567   | 3                     | 9.86  | 1106570 | 20.0 |
| Heptadecane          | 10.30 | 4.8     | ng    | 265947   | 3                     | 9.86  | 1106570 | 20.0 |
| Octacosane           | 10.52 | 4.2     | ng    | 229698   | 3                     | 9.86  | 1106570 | 20.0 |
| n-Hexadecanoic acid  | 11.88 | 4.1     | ng    | 208667   | 4                     | 11.35 | 1022610 | 20.0 |
| Pentadecanoic acid   | 12.66 | 2.6     | ng    | 135707   | 4                     | 11.35 | 1022610 | 20.0 |
| 1-Tetracosanol       | 13.83 | 5.1     | ng    | 166802   | 5                     | 13.98 | 654864  | 20.0 |