

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\  
 Data File : BM019139.D  
 Acq On : 13 Mar 2019 09:04  
 Operator : JU/SJ  
 Sample : K1824-17  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 P-3

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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\8270-BM031119.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         | 5.257       | 363           | 369         | 382          | rBV      | 3933550        | 5965979       | 40.64%          | 6.625%        |
| 2         | 6.846       | 631           | 639         | 657          | rBV      | 4341841        | 7064611       | 48.12%          | 7.845%        |
| 3         | 7.199       | 691           | 699         | 715          | rBV      | 4373903        | 6940448       | 47.27%          | 7.707%        |
| 4         | 7.657       | 771           | 777         | 787          | rBV      | 757487         | 1232264       | 8.39%           | 1.368%        |
| 5         | 7.975       | 823           | 831         | 841          | rBV      | 2925215        | 4656887       | 31.72%          | 5.171%        |
| 6         | 8.828       | 968           | 976         | 998          | rBV      | 2298404        | 4013116       | 27.34%          | 4.456%        |
| 7         | 10.445      | 1244          | 1251        | 1274         | rBV      | 969369         | 1759940       | 11.99%          | 1.954%        |
| 8         | 12.928      | 1666          | 1673        | 1695         | rBV      | 6115115        | 8948990       | 60.96%          | 9.938%        |
| 9         | 13.780      | 1813          | 1818        | 1829         | rBV      | 386865         | 627074        | 4.27%           | 0.696%        |
| 10        | 14.316      | 1899          | 1909        | 1928         | rBV2     | 1701980        | 2770469       | 18.87%          | 3.076%        |
| 11        | 15.816      | 2157          | 2164        | 2182         | rBV      | 6676467        | 9572159       | 65.20%          | 10.630%       |
| 12        | 17.068      | 2371          | 2377        | 2389         | rBV2     | 2368764        | 3555703       | 24.22%          | 3.948%        |
| 13        | 17.957      | 2523          | 2528        | 2540         | rBV      | 325098         | 554841        | 3.78%           | 0.616%        |
| 14        | 19.704      | 2819          | 2825        | 2839         | rBV      | 11667647       | 14681019      | 100.00%         | 16.303%       |
| 15        | 20.962      | 3035          | 3039        | 3052         | rBV      | 1531940        | 1855572       | 12.64%          | 2.061%        |
| 16        | 21.268      | 3085          | 3091        | 3101         | rBV      | 3121071        | 4186798       | 28.52%          | 4.649%        |
| 17        | 21.403      | 3111          | 3114        | 3119         | rBV2     | 119688         | 150350        | 1.02%           | 0.167%        |
| 18        | 21.833      | 3184          | 3187        | 3189         | rBV2     | 133664         | 170792        | 1.16%           | 0.190%        |
| 19        | 22.292      | 3261          | 3265        | 3269         | rBV2     | 208460         | 258560        | 1.76%           | 0.287%        |
| 20        | 22.792      | 3346          | 3350        | 3354         | rVB      | 304502         | 411576        | 2.80%           | 0.457%        |
| 21        | 23.356      | 3441          | 3446        | 3452         | rVB      | 476519         | 719426        | 4.90%           | 0.799%        |
| 22        | 23.503      | 3464          | 3471        | 3482         | rBV      | 1853428        | 3555077       | 24.22%          | 3.948%        |
| 23        | 23.997      | 3549          | 3555        | 3563         | rVB      | 663511         | 1276552       | 8.70%           | 1.418%        |
| 24        | 24.086      | 3566          | 3570        | 3581         | rVB3     | 295048         | 602016        | 4.10%           | 0.669%        |
| 25        | 24.739      | 3675          | 3681        | 3688         | rVB      | 785150         | 1608910       | 10.96%          | 1.787%        |
| 26        | 25.603      | 3821          | 3828        | 3837         | rVB      | 627207         | 1585298       | 10.80%          | 1.760%        |
| 27        | 26.621      | 3995          | 4001        | 4010         | rVB2     | 481315         | 1328287       | 9.05%           | 1.475%        |

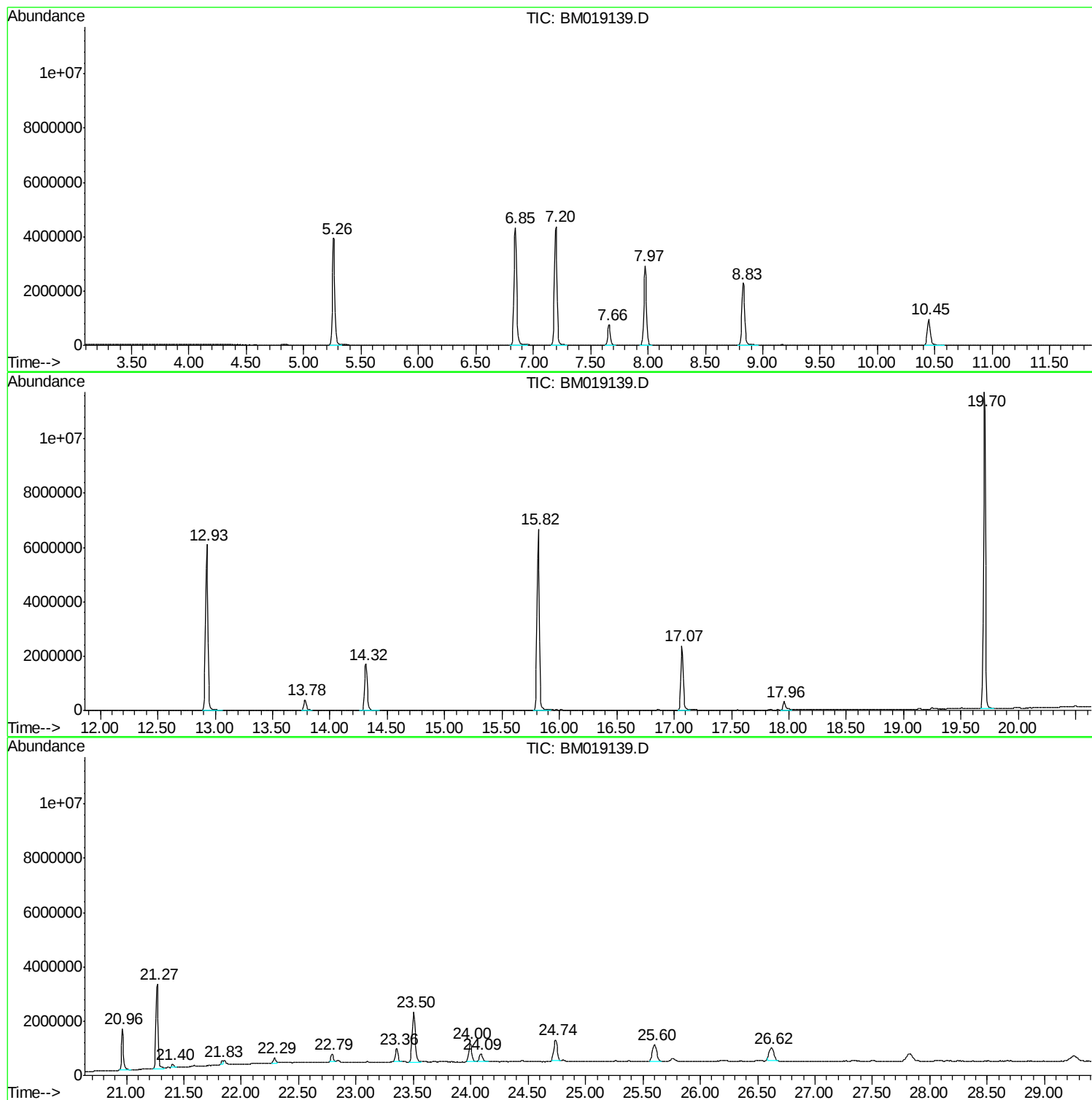
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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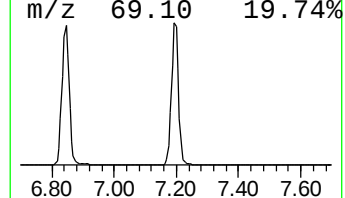
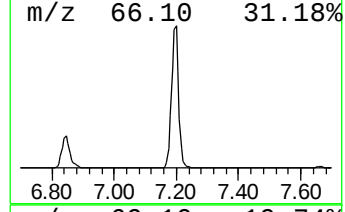
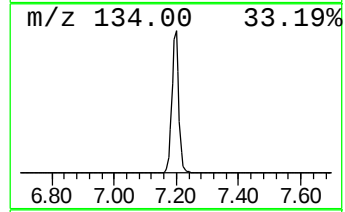
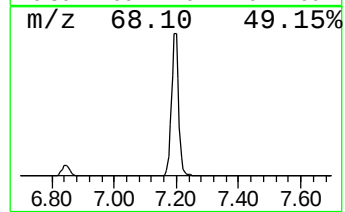
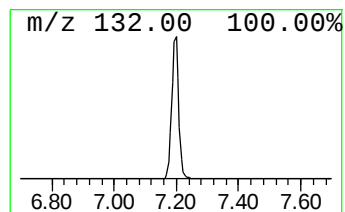
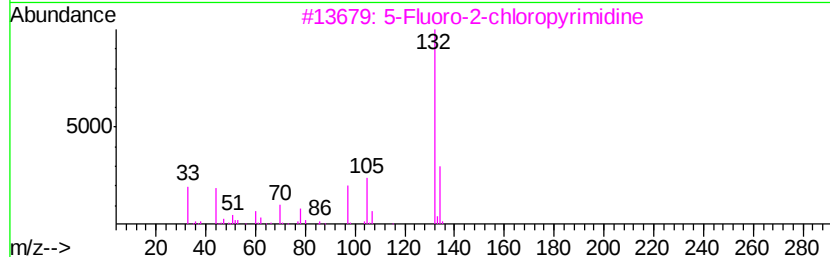
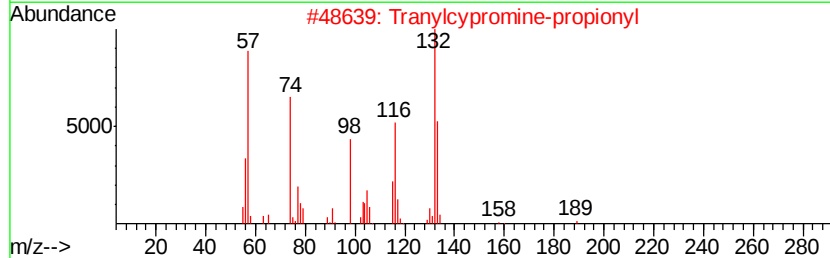
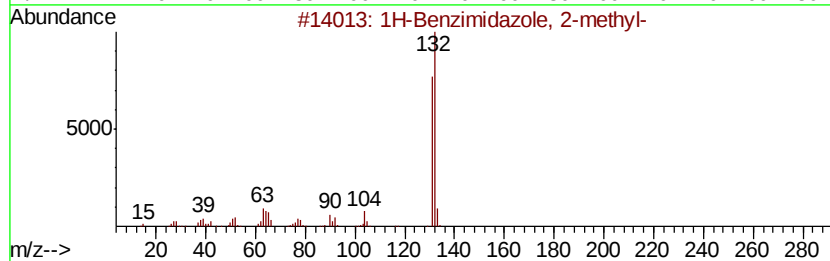
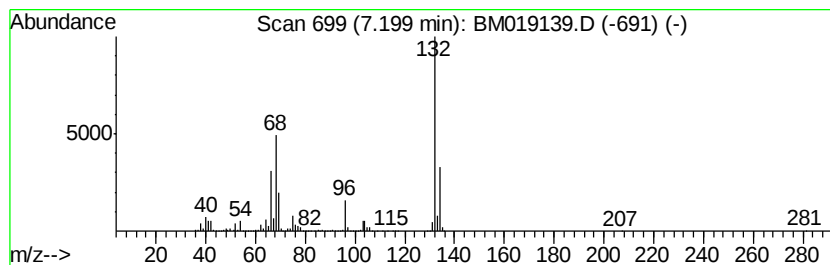
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown7.20 Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 7.20 | 112.65 ng | 6940450 | 1,4-Dichlorobenzene-d4 | 7.66 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 22   |
| 2    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 3    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 12   |
| 4    |    | Benzene, 1,2,4-trifluoro-        | 132 | C6H3F3    | 000367-23-7  | 9    |
| 5    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 9    |



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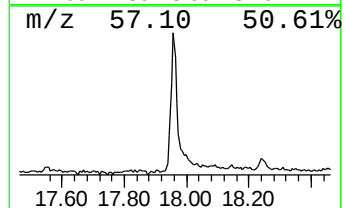
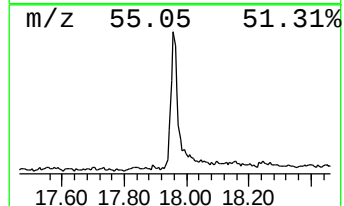
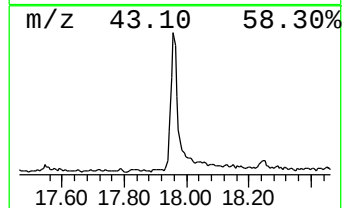
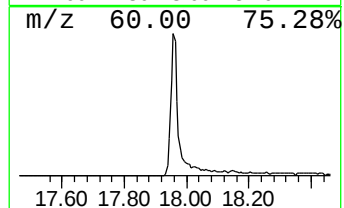
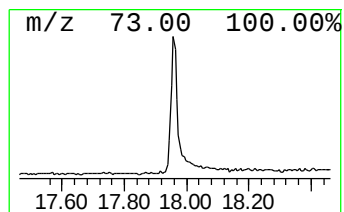
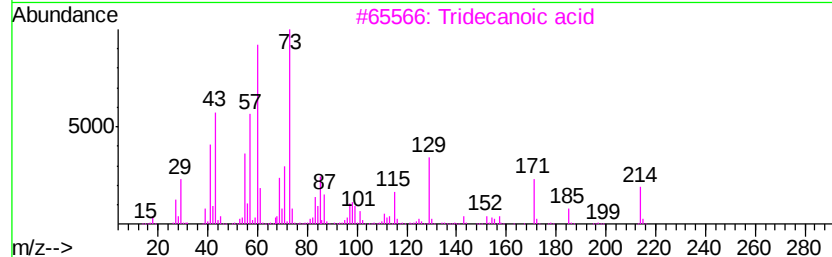
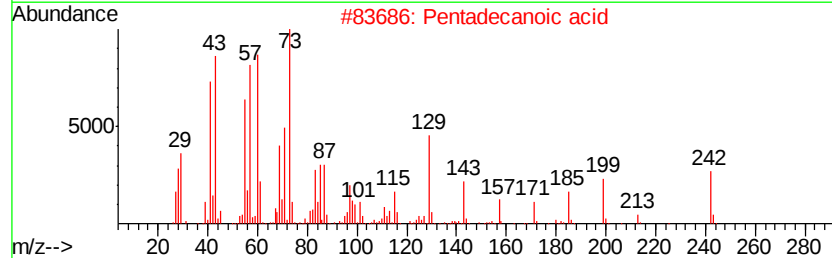
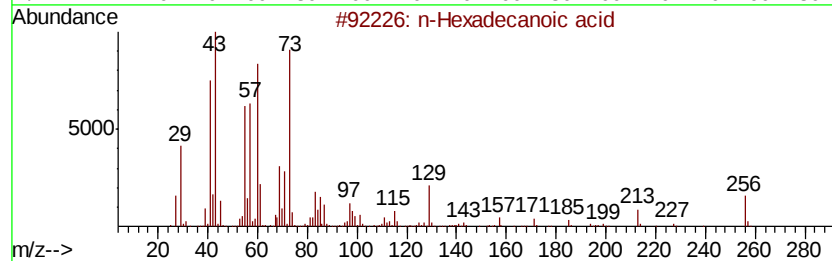
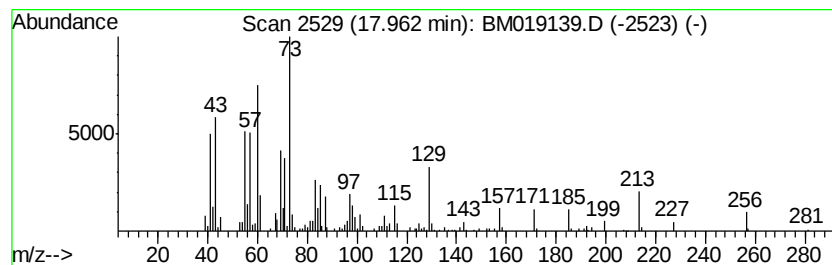
Instrument :  
BNA\_M  
ClientSampleId :  
P-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 n-Hexadecanoic acid Concentration Rank 10

| R.T.    | EstConc | Area                | Relative to ISTD | R.T.           |
|---------|---------|---------------------|------------------|----------------|
| 17.96   | 3.12 ng | 554841              | Phenanthrene-d10 | 17.07          |
| Hit# of | 5       | Tentative ID        | MW MolForm       | CAS# Qual      |
| 1       |         | n-Hexadecanoic acid | 256 C16H32O2     | 000057-10-3 98 |
| 2       |         | Pentadecanoic acid  | 242 C15H30O2     | 001002-84-2 72 |
| 3       |         | Tridecanoic acid    | 214 C13H26O2     | 000638-53-9 59 |
| 4       |         | Tetradecanoic acid  | 228 C14H28O2     | 000544-63-8 47 |
| 5       |         | n-Decanoic acid     | 172 C10H20O2     | 000334-48-5 43 |



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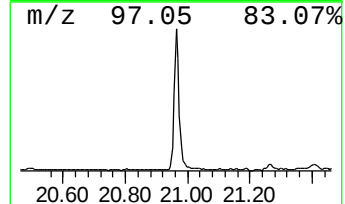
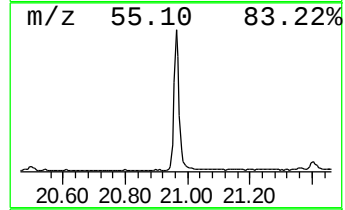
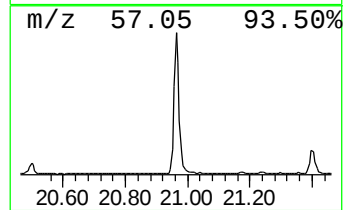
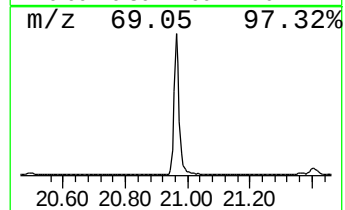
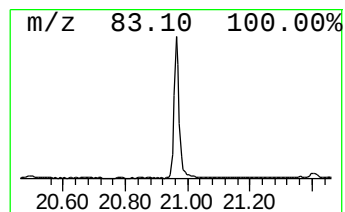
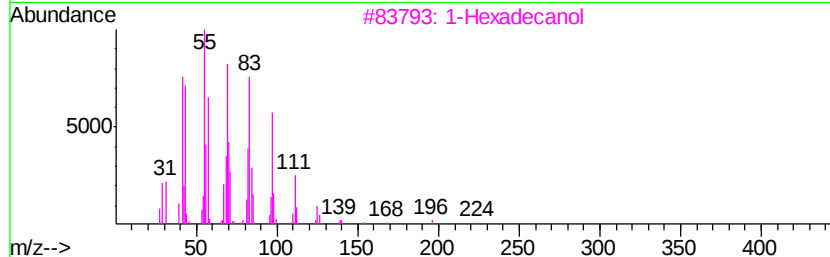
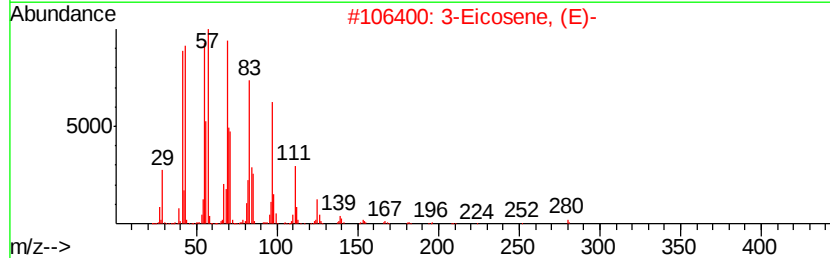
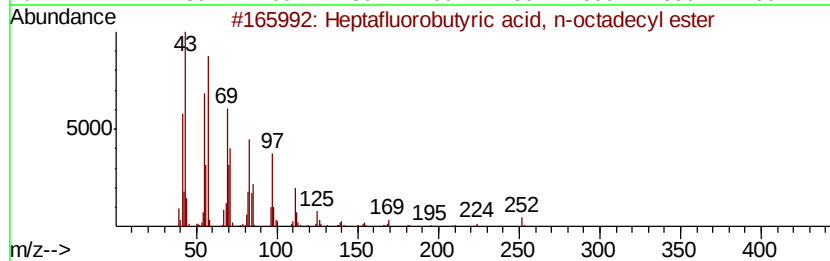
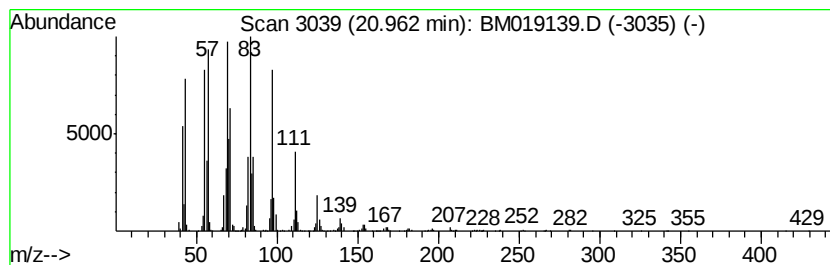
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Heptafluorobutyric acid, n-... Concentration Rank 5

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 20.96 | 8.86 ng | 1855570 | Chrysene-d12     | 21.27 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|-----------|-------------------------------------|-----|------------|-------------|------|
| 1         | Heptafluorobutyric acid, n-octad... | 466 | C22H37F7O2 | 000400-57-7 | 91   |
| 2         | 3-Eicosene, (E)-                    | 280 | C20H40     | 074685-33-9 | 91   |
| 3         | 1-Hexadecanol                       | 242 | C16H34O    | 036653-82-4 | 90   |
| 4         | 1-Tetracosanol                      | 354 | C24H50O    | 000506-51-4 | 87   |
| 5         | 9-Eicosene, (E)-                    | 280 | C20H40     | 074685-29-3 | 83   |



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P-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M  
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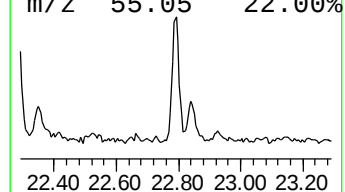
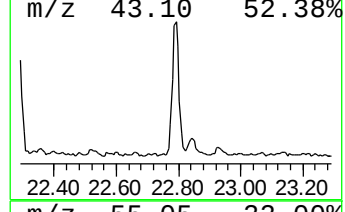
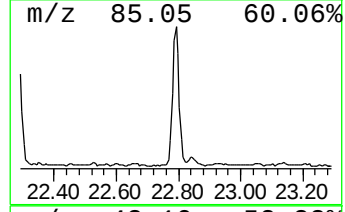
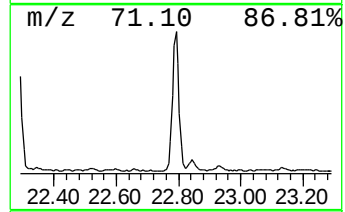
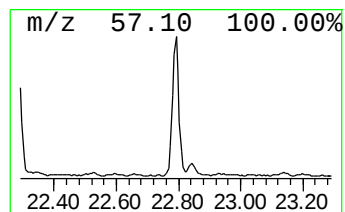
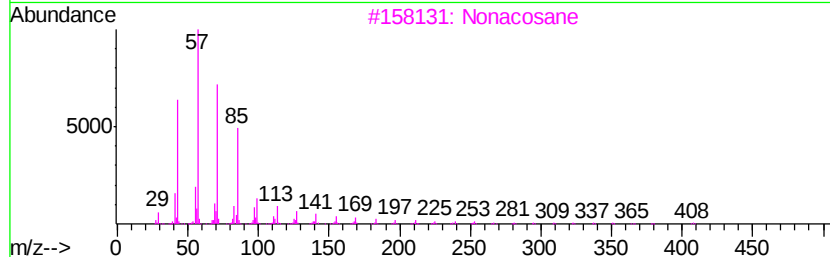
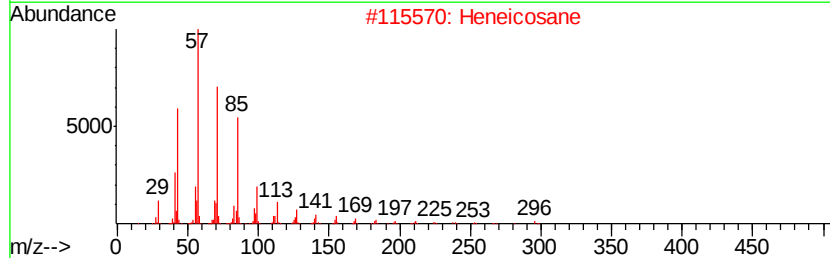
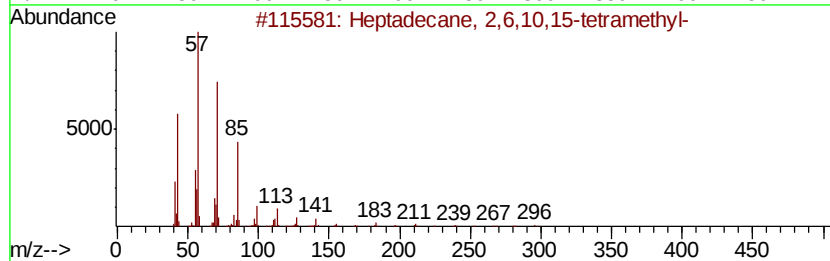
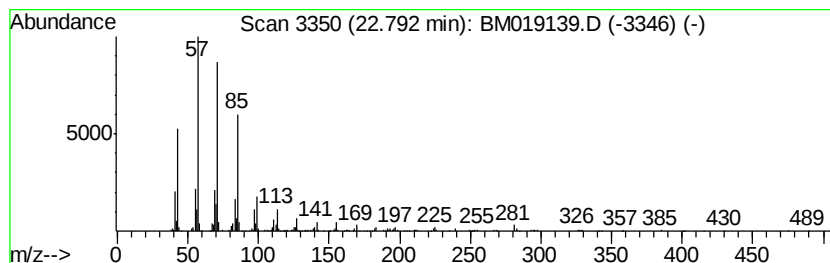
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Peak Number 5 Heptadecane, 2,6,10,15-tetr... Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 22.79 | 2.32 ng | 411576 | Perylene-d12     | 23.50 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 93   |
| 2    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    |   | Nonacosane                          | 408 | C29H60  | 000630-03-5 | 91   |
| 4    |    |   | Heptadecane, 9-octyl-               | 352 | C25H52  | 007225-64-1 | 90   |
| 5    |    |   | Docosane                            | 310 | C22H46  | 000629-97-0 | 90   |



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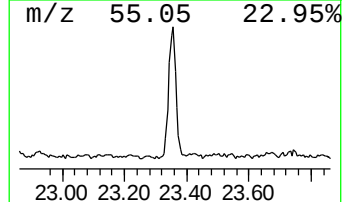
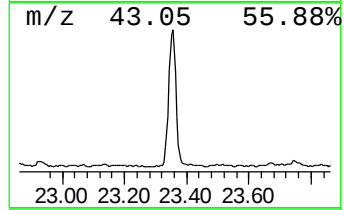
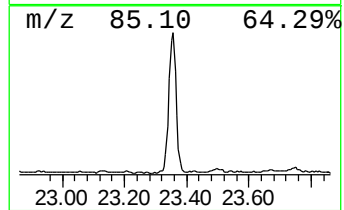
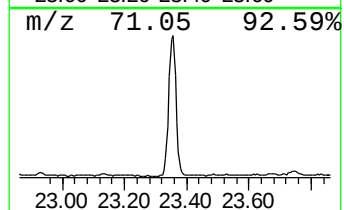
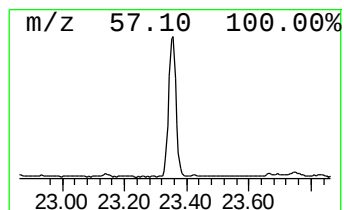
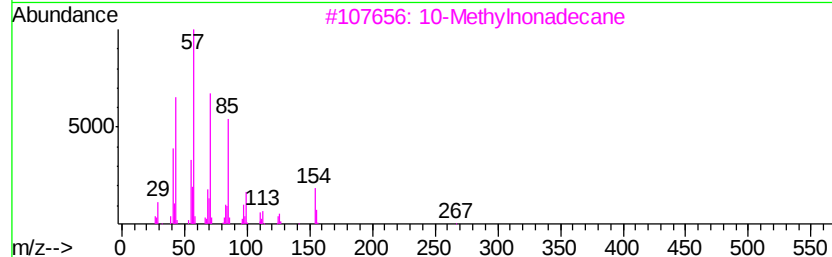
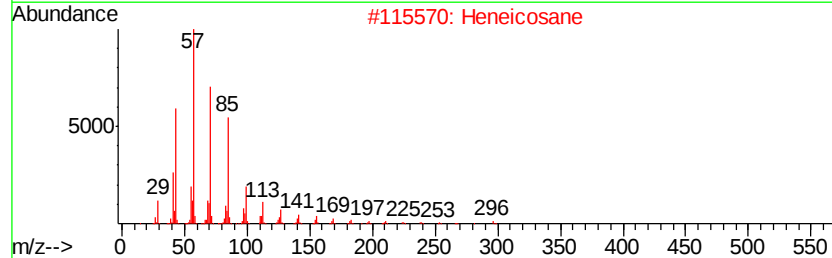
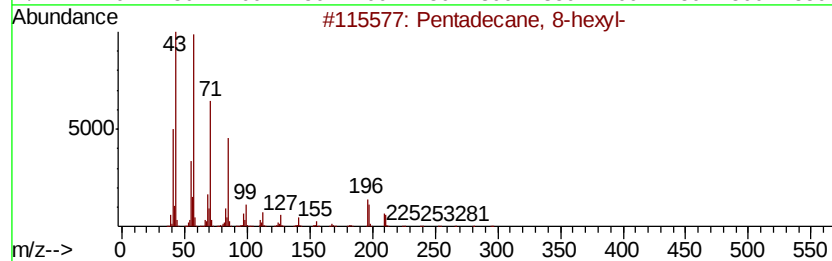
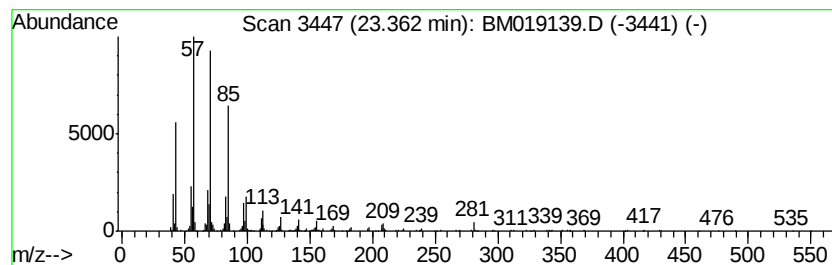
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 6 Pentadecane, 8-hexyl- Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 23.36 | 4.05 ng | 719426 | Perylene-d12     | 23.50 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 90   |
| 2    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 87   |
| 3    |    | 10-Methylnonadecane   | 282 | C20H42  | 056862-62-5 | 86   |
| 4    |    | Octadecane, 1-iodo-   | 380 | C18H37I | 000629-93-6 | 86   |
| 5    |    | Tridecane, 6-propyl-  | 226 | C16H34  | 055045-10-8 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\  
Data File : BM019139.D  
Acq On : 13 Mar 2019 09:04  
Operator : JU/SJ  
Sample : K1824-17  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
P-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

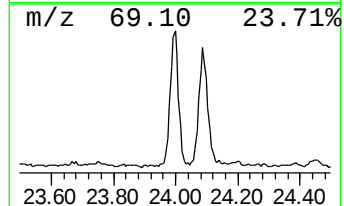
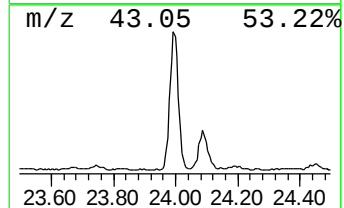
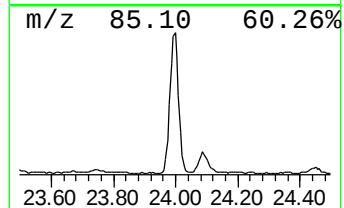
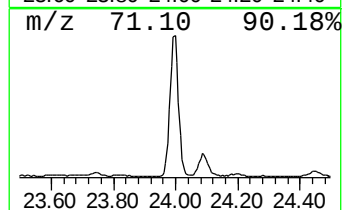
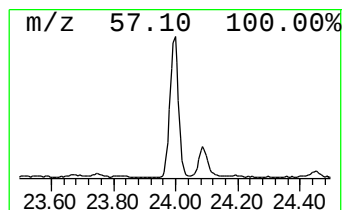
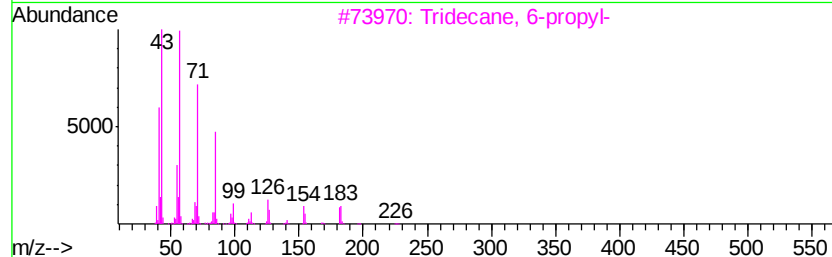
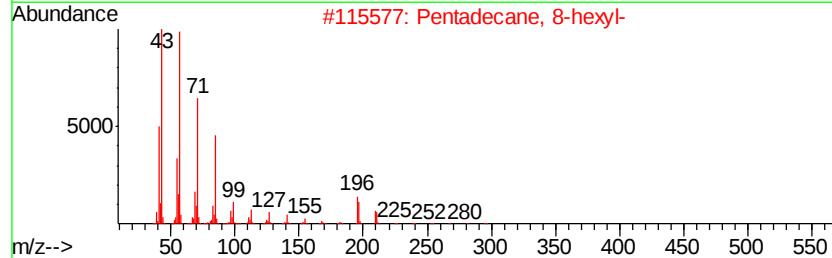
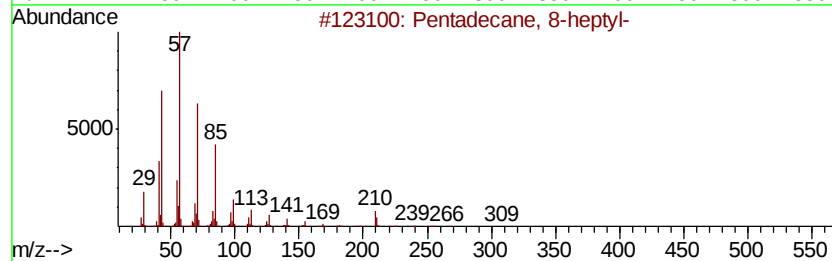
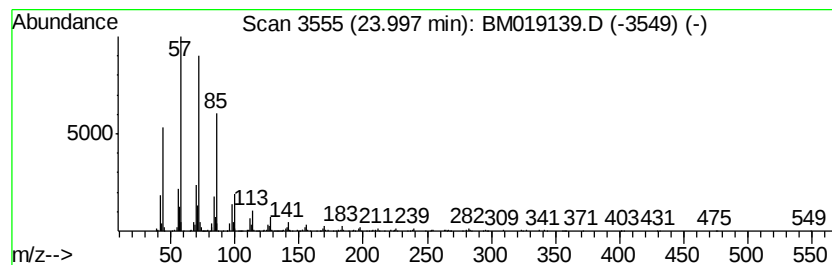
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Pentadecane, 8-heptyl- Concentration Rank 7

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 24.00 | 7.18 ng | 1276550 | Perylene-d12     | 23.50 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 8-heptyl-           | 310 | C22H46  | 071005-15-7 | 91   |
| 2    |    | Pentadecane, 8-hexyl-            | 296 | C21H44  | 013475-75-7 | 90   |
| 3    |    | Tridecane, 6-propyl-             | 226 | C16H34  | 055045-10-8 | 87   |
| 4    |    | Octadecane                       | 254 | C18H38  | 000593-45-3 | 87   |
| 5    |    | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\  
Data File : BM019139.D  
Acq On : 13 Mar 2019 09:04  
Operator : JU/SJ  
Sample : K1824-17  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
P-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

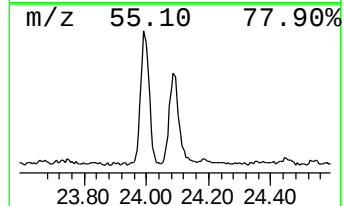
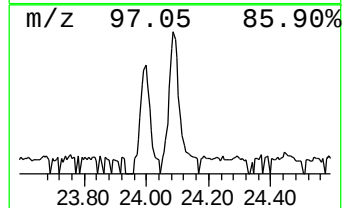
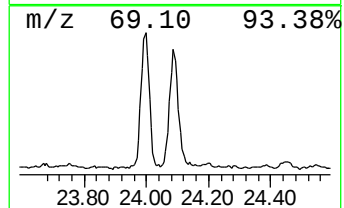
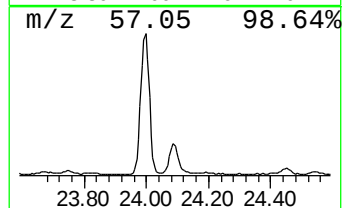
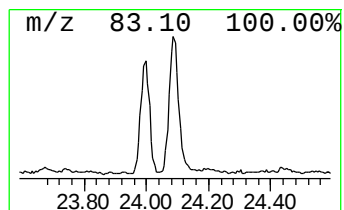
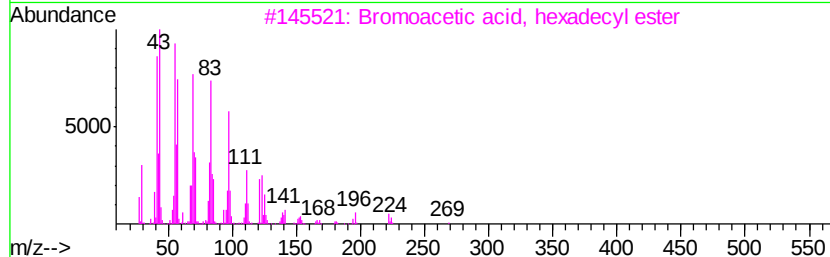
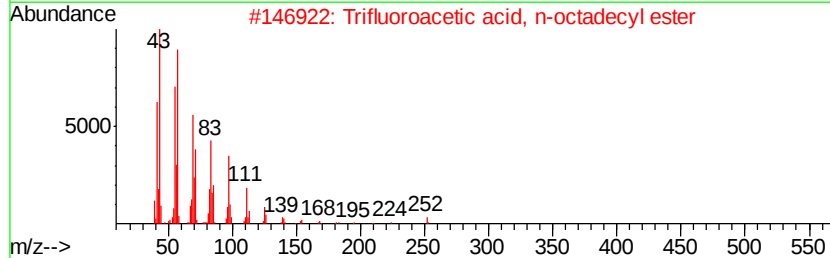
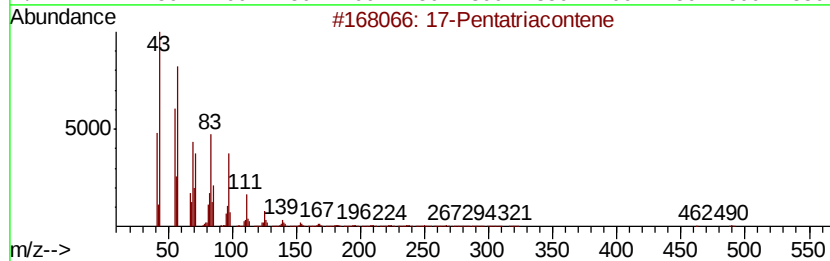
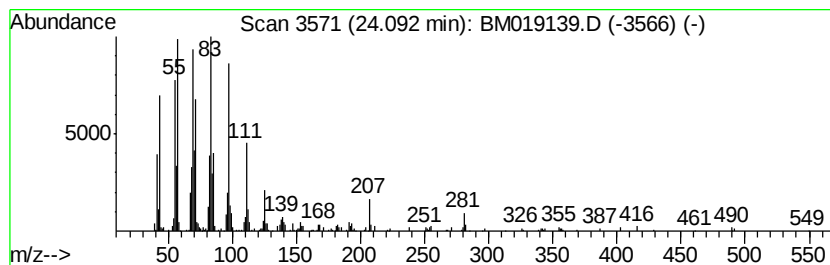
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 17-Pentatriacontene Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 24.09 | 3.39 ng | 602016 | Perylene-d12     | 23.50 |

| Hit# of | 5                                    | Tentative ID | MW  | MolForm    | CAS#        | Qual |
|---------|--------------------------------------|--------------|-----|------------|-------------|------|
| 1       | 17-Pentatriacontene                  |              | 491 | C35H70     | 006971-40-0 | 93   |
| 2       | Trifluoroacetic acid, n-octadecyl... |              | 366 | C20H37F3O2 | 079392-43-1 | 91   |
| 3       | Bromoacetic acid, hexadecyl ester    |              | 362 | C18H35BrO2 | 005454-48-8 | 87   |
| 4       | 1-Heneicosyl formate                 |              | 340 | C22H44O2   | 077899-03-7 | 87   |
| 5       | 3-Eicosene, (E)-                     |              | 280 | C20H40     | 074685-33-9 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\  
Data File : BM019139.D  
Acq On : 13 Mar 2019 09:04  
Operator : JU/SJ  
Sample : K1824-17  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
P-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

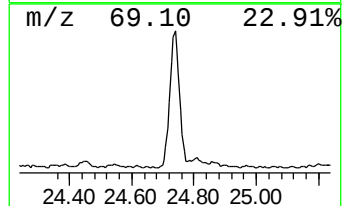
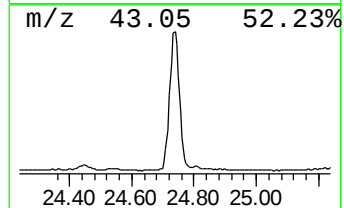
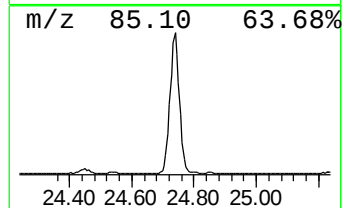
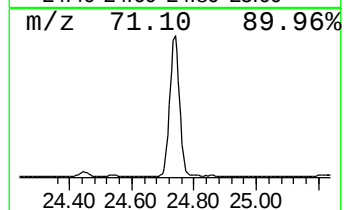
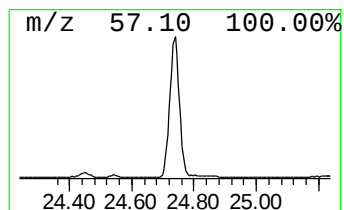
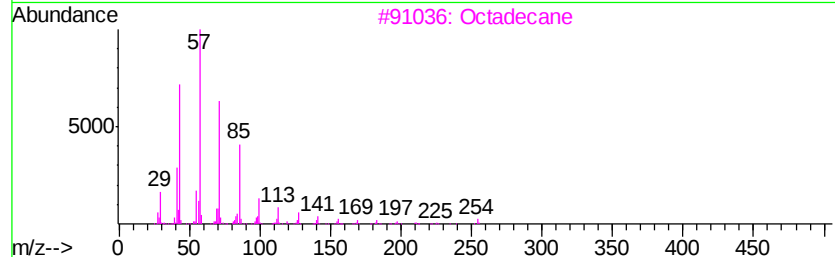
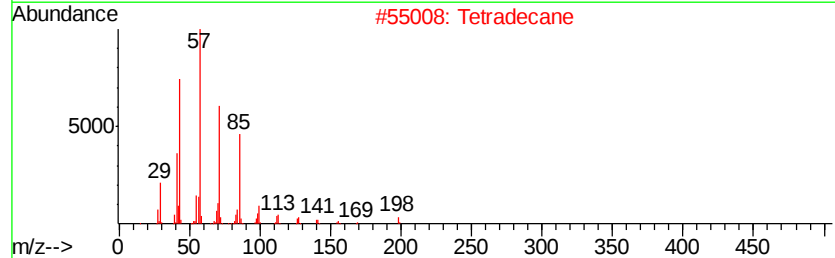
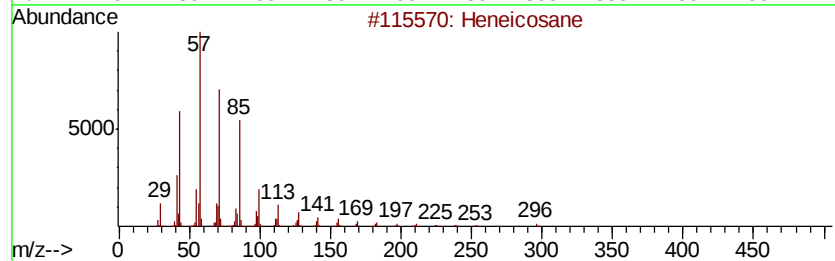
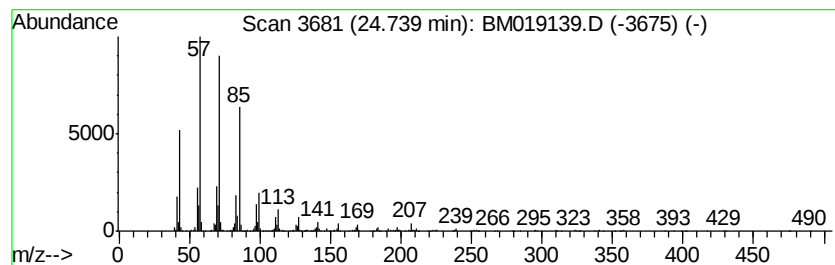
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Heneicosane Concentration Rank 3

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 24.74 | 9.05 ng | 1608910 | Perylene-d12     | 23.50 |

| Hit# of | 5                     | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-----------------------|--------------|-----|---------|-------------|------|
| 1       | Heneicosane           |              | 296 | C21H44  | 000629-94-7 | 91   |
| 2       | Tetradecane           |              | 198 | C14H30  | 000629-59-4 | 87   |
| 3       | Octadecane            |              | 254 | C18H38  | 000593-45-3 | 87   |
| 4       | Heptacosane           |              | 380 | C27H56  | 000593-49-7 | 87   |
| 5       | Pentadecane, 8-hexyl- |              | 296 | C21H44  | 013475-75-7 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\  
Data File : BM019139.D  
Acq On : 13 Mar 2019 09:04  
Operator : JU/SJ  
Sample : K1824-17  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
P-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

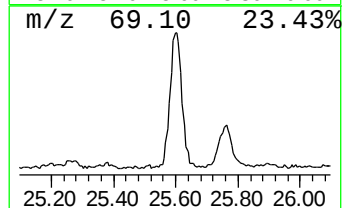
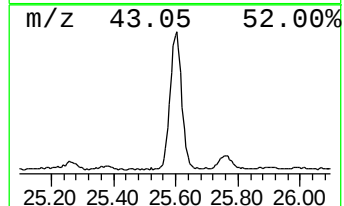
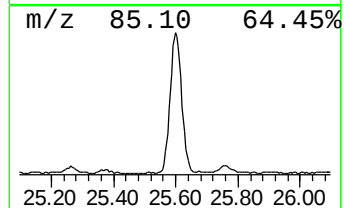
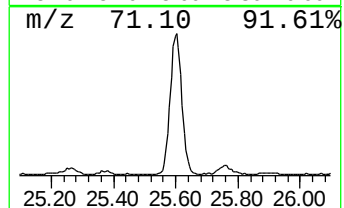
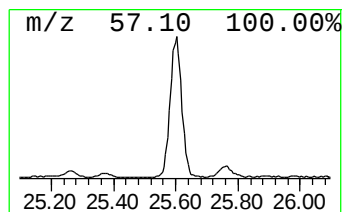
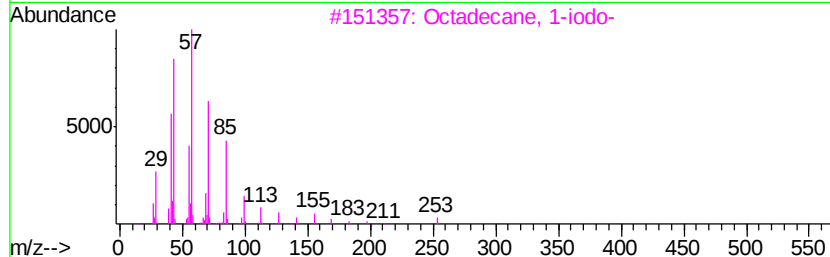
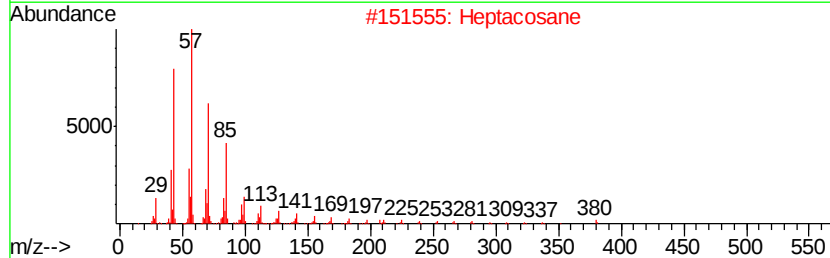
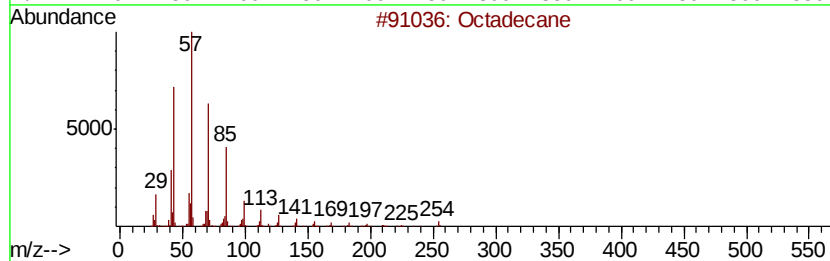
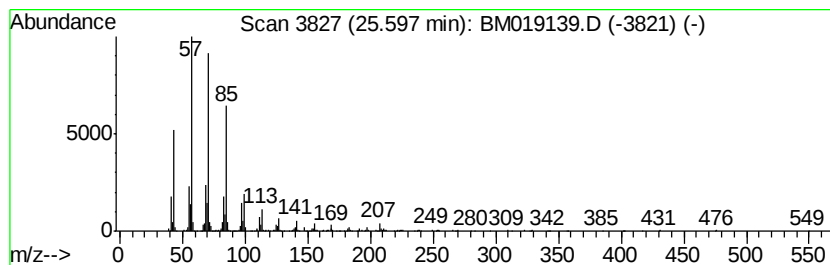
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Octadecane Concentration Rank 4

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 25.60 | 8.92 ng | 1585300 | Perylene-d12     | 23.50 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Octadecane             | 254 | C18H38  | 000593-45-3 | 87   |
| 2    |    |   | Heptacosane            | 380 | C27H56  | 000593-49-7 | 87   |
| 3    |    |   | Octadecane, 1-iodo-    | 380 | C18H37I | 000629-93-6 | 87   |
| 4    |    |   | Docosane, 5-butyl-     | 366 | C26H54  | 055282-16-1 | 86   |
| 5    |    |   | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031219\  
Data File : BM019139.D  
Acq On : 13 Mar 2019 09:04  
Operator : JU/SJ  
Sample : K1824-17  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
P-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

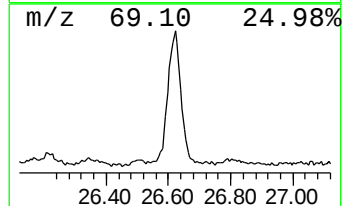
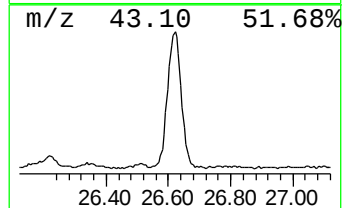
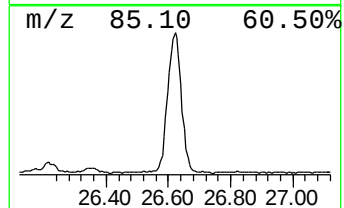
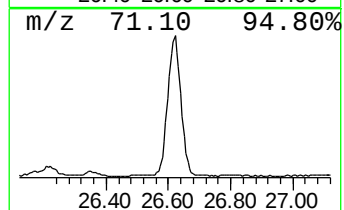
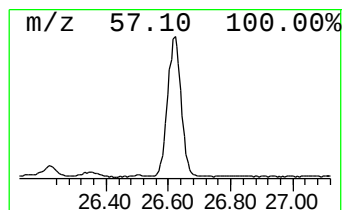
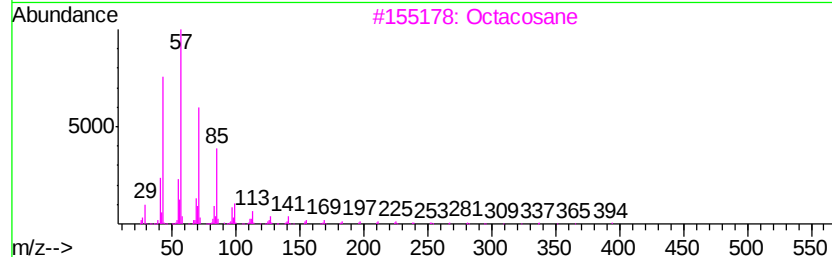
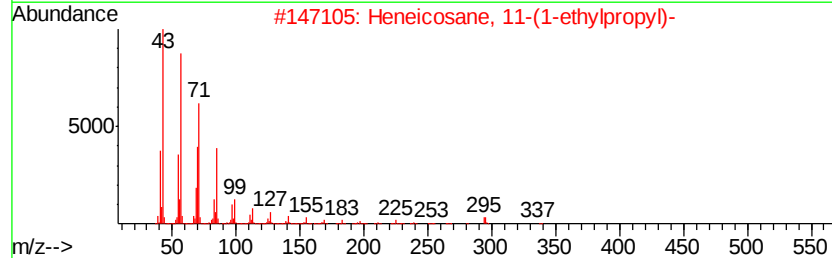
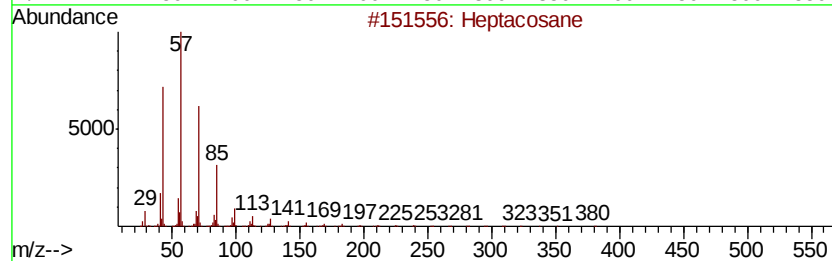
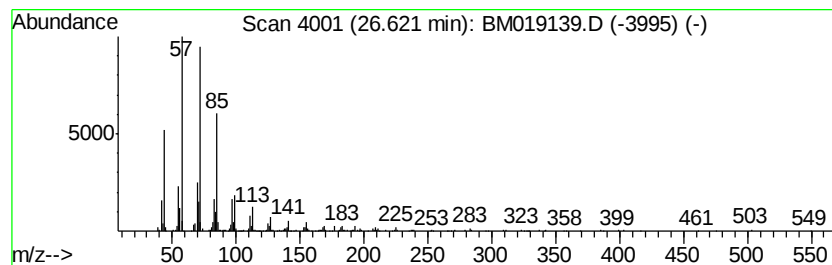
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Heptacosane Concentration Rank 6

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 26.62 | 7.47 ng | 1328290 | Perylene-d12     | 23.50 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptacosane                      | 380 | C27H56  | 000593-49-7 | 91   |
| 2    |    | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 91   |
| 3    |    | Octacosane                       | 394 | C28H58  | 000630-02-4 | 87   |
| 4    |    | Tridecane, 6-propyl-             | 226 | C16H34  | 055045-10-8 | 87   |
| 5    |    | Pentadecane, 8-heptyl-           | 310 | C22H46  | 071005-15-7 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM031219\  
Data File : BM019139.D  
Acq On : 13 Mar 2019 09:04  
Operator : JU/SJ  
Sample : K1824-17  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
P-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\8270-BM031119.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 |
| unknown7.20          | 7.20  | 112.7   | ng    | 6940450  | 1                     | 7.66  | 1232260 | 20.0 |
| n-Hexadecanoic acid  | 17.96 | 3.1     | ng    | 554841   | 4                     | 17.07 | 3555700 | 20.0 |
| Heptafluorobutyri... | 20.96 | 8.9     | ng    | 1855570  | 5                     | 21.27 | 4186800 | 20.0 |
| Heptadecane, 2,6,... | 22.79 | 2.3     | ng    | 411576   | 6                     | 23.50 | 3555080 | 20.0 |
| Pentadecane, 8-he... | 23.36 | 4.0     | ng    | 719426   | 6                     | 23.50 | 3555080 | 20.0 |
| Pentadecane, 8-he... | 24.00 | 7.2     | ng    | 1276550  | 6                     | 23.50 | 3555080 | 20.0 |
| 17-Pentatriacontene  | 24.09 | 3.4     | ng    | 602016   | 6                     | 23.50 | 3555080 | 20.0 |
| Heneicosane          | 24.74 | 9.1     | ng    | 1608910  | 6                     | 23.50 | 3555080 | 20.0 |
| Octadecane           | 25.60 | 8.9     | ng    | 1585300  | 6                     | 23.50 | 3555080 | 20.0 |
| Heptacosane          | 26.62 | 7.5     | ng    | 1328290  | 6                     | 23.50 | 3555080 | 20.0 |