

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\  
 Data File : BF109903.D  
 Acq On : 16 Oct 2018 14:20  
 Operator : JU/SJ  
 Sample : J5515-10  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 AOC3-01A

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\_F\METHODS\8270-BF101518.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         | 2.122       | 3             | 6           | 12           | rVB      | 358602         | 520518        | 10.02%          | 1.065%        |
| 2         | 2.328       | 34            | 41          | 49           | rVB      | 701060         | 912034        | 17.56%          | 1.867%        |
| 3         | 3.587       | 253           | 255         | 270          | rBV      | 71363          | 141533        | 2.73%           | 0.290%        |
| 4         | 5.222       | 528           | 533         | 538          | rVB      | 166676         | 195942        | 3.77%           | 0.401%        |
| 5         | 5.604       | 593           | 598         | 601          | rBV      | 4983193        | 4818524       | 92.79%          | 9.862%        |
| 6         | 6.598       | 762           | 767         | 771          | rBV      | 4911205        | 5193166       | 100.00%         | 10.629%       |
| 7         | 6.751       | 788           | 793         | 796          | rBV      | 5041965        | 4962653       | 95.56%          | 10.157%       |
| 8         | 6.981       | 828           | 832         | 835          | rBV      | 1904663        | 1472894       | 28.36%          | 3.015%        |
| 9         | 7.139       | 854           | 859         | 862          | rBV      | 4132604        | 3898295       | 75.07%          | 7.979%        |
| 10        | 7.539       | 922           | 927         | 930          | rBV      | 3458577        | 3129125       | 60.25%          | 6.404%        |
| 11        | 8.263       | 1046          | 1050        | 1053         | rBV      | 2214334        | 1827809       | 35.20%          | 3.741%        |
| 12        | 9.339       | 1228          | 1233        | 1236         | rVB      | 5277220        | 5051941       | 97.28%          | 10.340%       |
| 13        | 9.728       | 1295          | 1299        | 1302         | rVB      | 825321         | 637184        | 12.27%          | 1.304%        |
| 14        | 10.022      | 1345          | 1349        | 1353         | rVB      | 1987601        | 1711931       | 32.97%          | 3.504%        |
| 15        | 10.810      | 1479          | 1483        | 1487         | rBV      | 2525346        | 2113014       | 40.69%          | 4.325%        |
| 16        | 11.510      | 1599          | 1602        | 1606         | rBV      | 1497969        | 1395541       | 26.87%          | 2.856%        |
| 17        | 12.016      | 1683          | 1688        | 1698         | rBV4     | 120587         | 205284        | 3.95%           | 0.420%        |
| 18        | 12.804      | 1817          | 1822        | 1833         | rBV      | 201258         | 348140        | 6.70%           | 0.713%        |
| 19        | 13.098      | 1868          | 1872        | 1875         | rBV      | 4131472        | 3699820       | 71.24%          | 7.572%        |
| 20        | 13.657      | 1964          | 1967        | 1970         | rVB      | 95643          | 74739         | 1.44%           | 0.153%        |
| 21        | 13.980      | 2018          | 2022        | 2026         | rBV2     | 383173         | 442579        | 8.52%           | 0.906%        |
| 22        | 14.157      | 2048          | 2052        | 2056         | rBV      | 1737115        | 1558608       | 30.01%          | 3.190%        |
| 23        | 14.304      | 2073          | 2077        | 2080         | rBV      | 248622         | 242376        | 4.67%           | 0.496%        |
| 24        | 14.610      | 2126          | 2129        | 2133         | rBV      | 335482         | 300646        | 5.79%           | 0.615%        |
| 25        | 14.933      | 2180          | 2184        | 2188         | rBV      | 360499         | 413273        | 7.96%           | 0.846%        |
| 26        | 15.010      | 2195          | 2197        | 2201         | rVB      | 101547         | 92100         | 1.77%           | 0.189%        |
| 27        | 15.286      | 2241          | 2244        | 2251         | rVB      | 430241         | 481809        | 9.28%           | 0.986%        |
| 28        | 15.692      | 2307          | 2313        | 2318         | rBV2     | 1497112        | 2115620       | 40.74%          | 4.330%        |
| 29        | 16.162      | 2388          | 2393        | 2398         | rVB      | 319577         | 492012        | 9.47%           | 1.007%        |
| 30        | 16.704      | 2481          | 2485        | 2492         | rVB      | 227687         | 410101        | 7.90%           | 0.839%        |

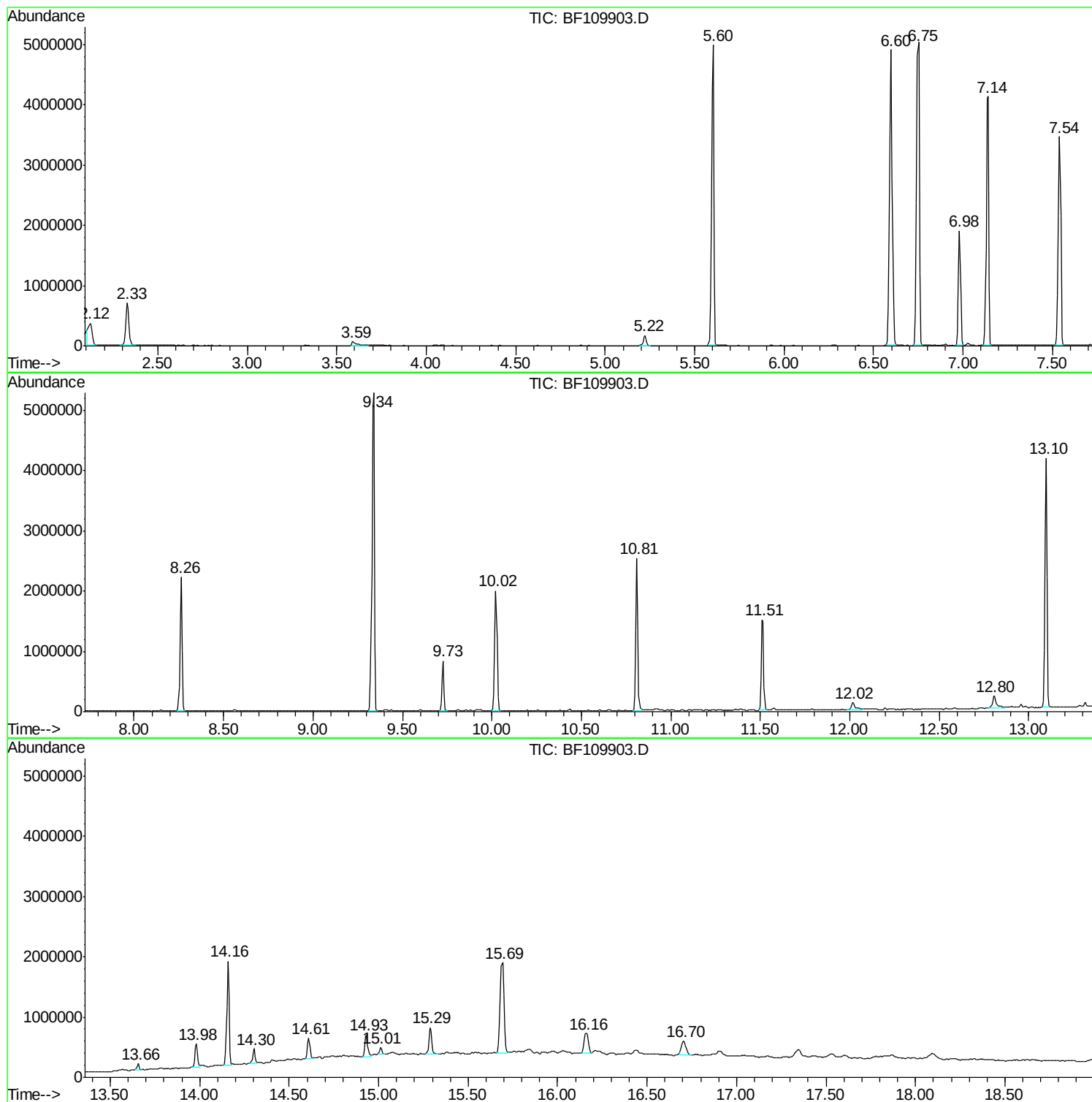
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P



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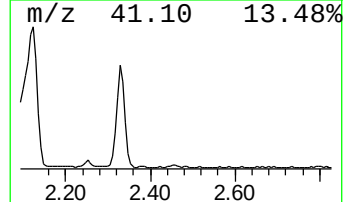
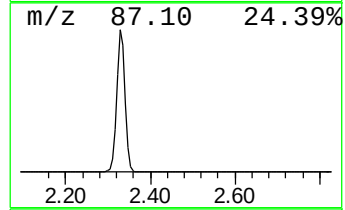
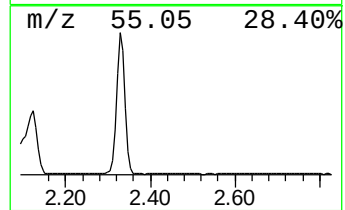
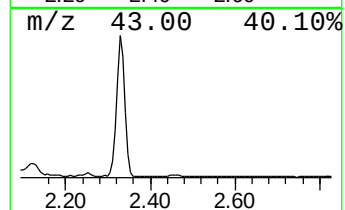
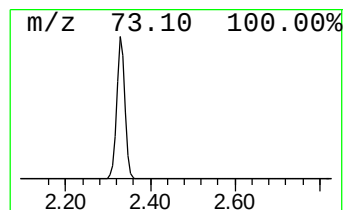
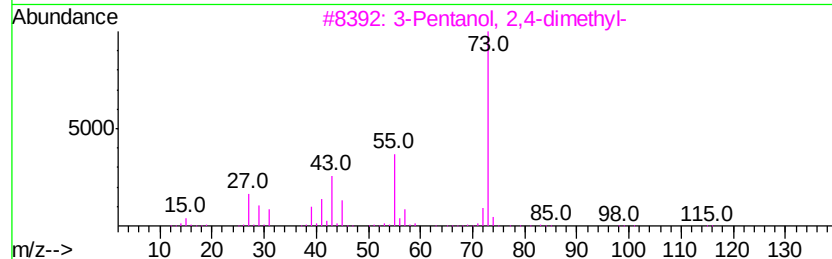
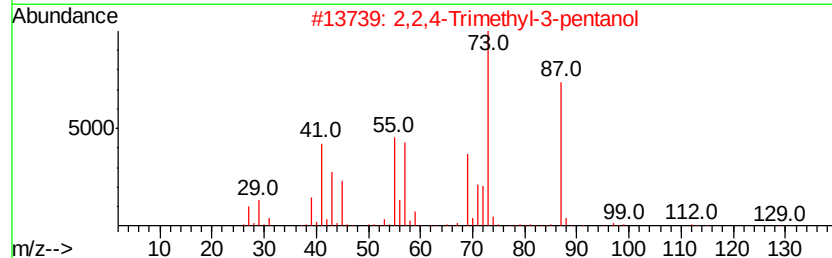
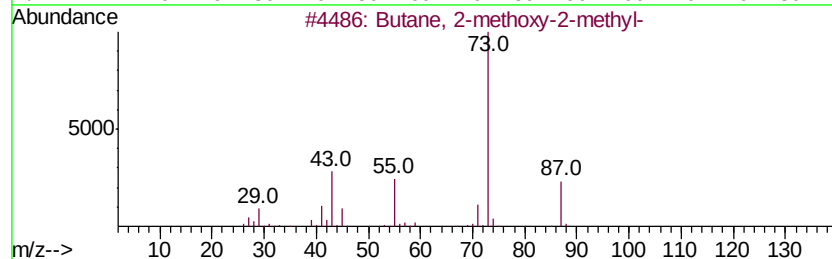
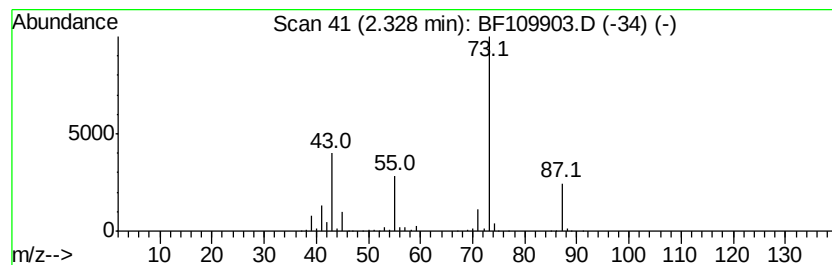
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 2.33 | 12.38 ng | 912034 | 1,4-Dichlorobenzene-d4 | 6.98 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 39   |
| 3    |    | 3-Pentanol, 2,4-dimethyl-   | 116 | C7H16O  | 000600-36-2 | 37   |
| 4    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 25   |
| 5    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 12   |



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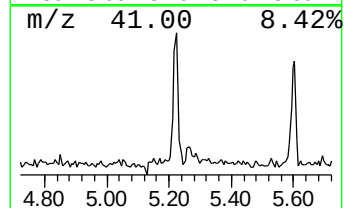
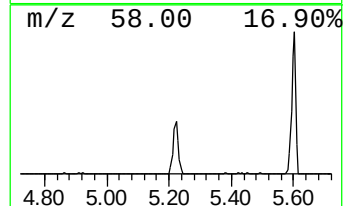
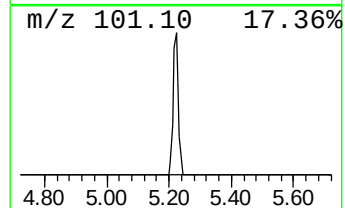
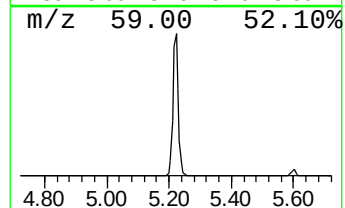
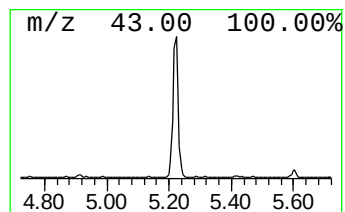
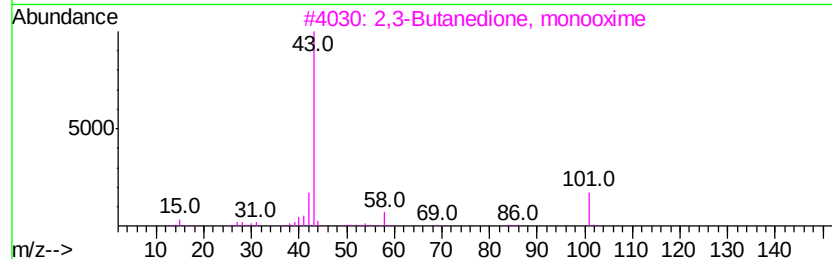
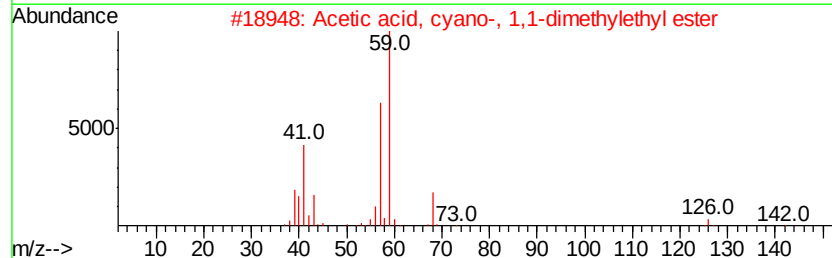
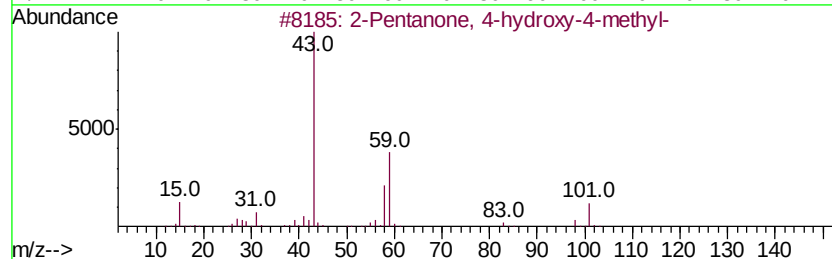
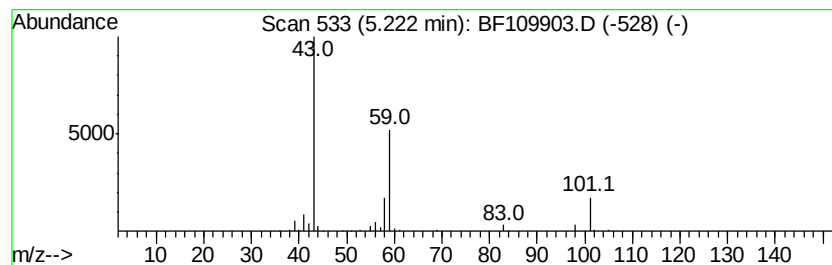
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.22 | 2.66 ng | 195942 | 1,4-Dichlorobenzene-d4 | 6.98 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 59   |
| 2    |    | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 3    |    | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 10   |
| 4    |    | 5-Hexen-2-one                        | 98  | C6H10O   | 000109-49-9 | 9    |
| 5    |    | Acetone                              | 58  | C3H6O    | 000067-64-1 | 9    |



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ALS Vial : 5 Sample Multiplier: 1

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ClientSampleId :  
AOC3-01A

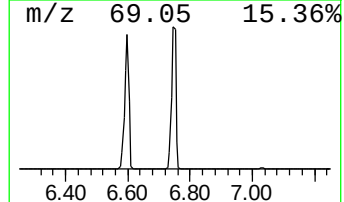
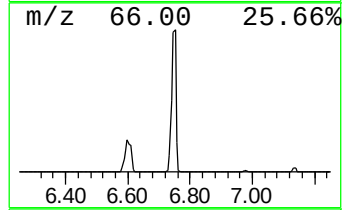
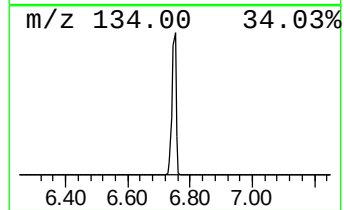
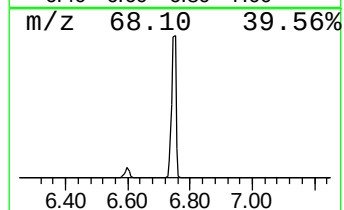
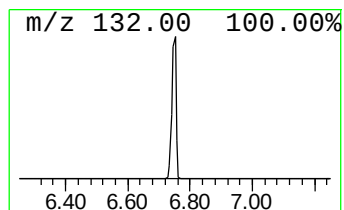
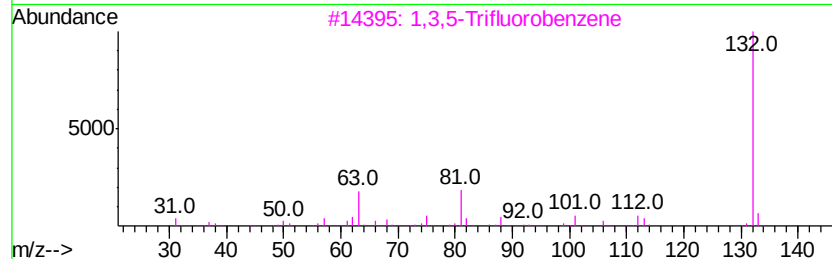
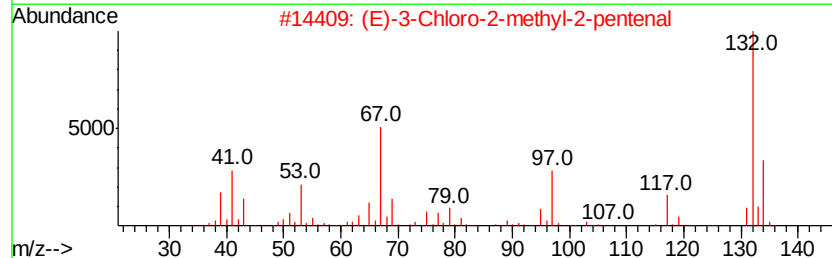
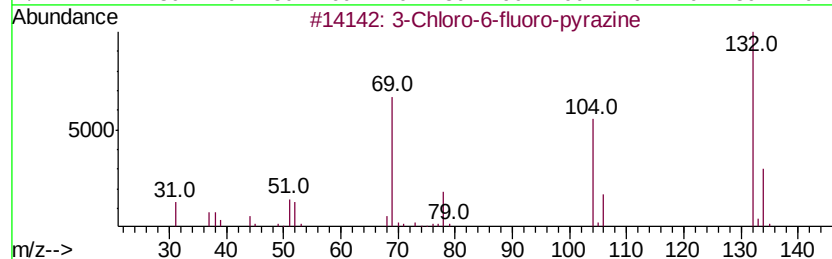
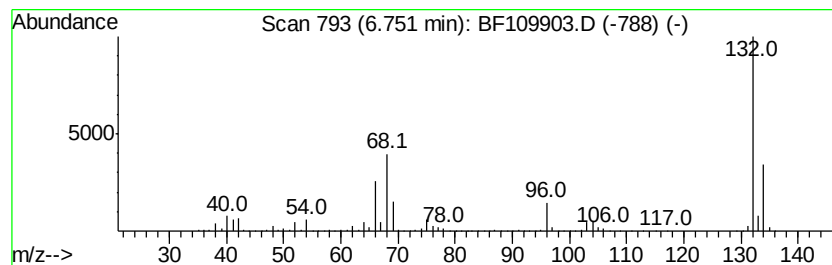
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown6.75 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.75 | 67.39 ng | 4962650 | 1,4-Dichlorobenzene-d4 | 6.98 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    |    | 1,3,5-Trifluorobenzene              | 132 | C6H3F3    | 000372-38-3  | 9    |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0  | 9    |



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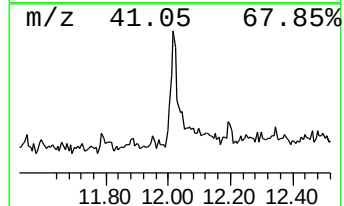
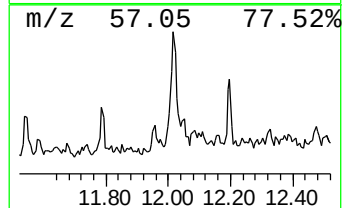
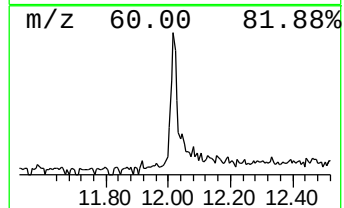
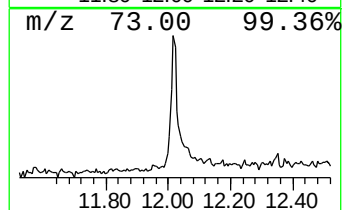
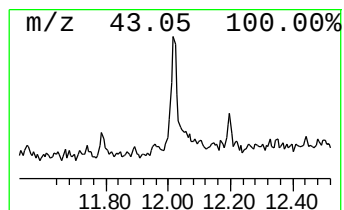
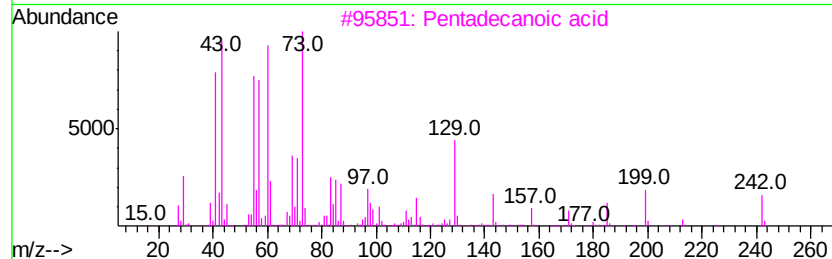
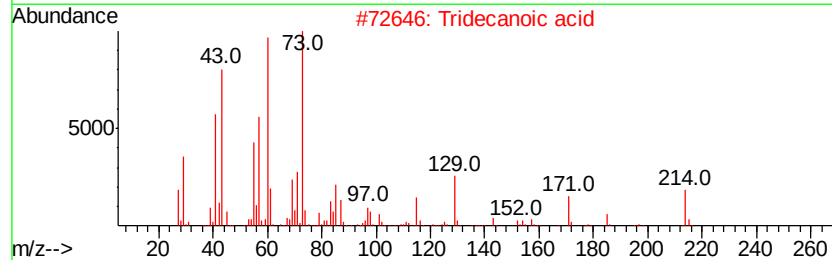
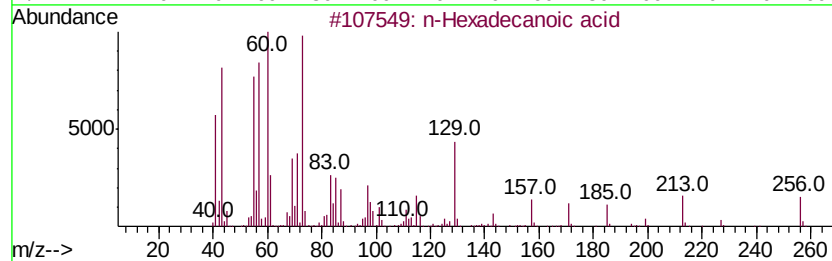
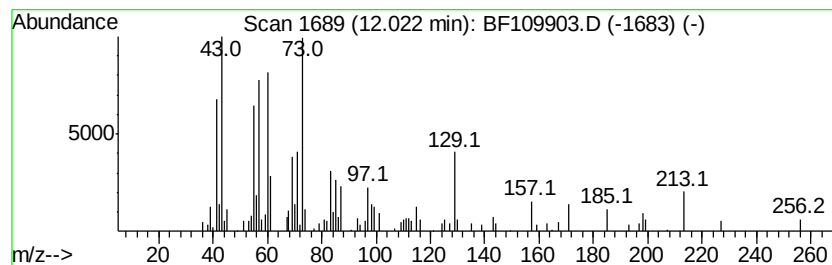
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 13

| R.T.    | EstConc | Area                | Relative to ISTD | R.T.           |
|---------|---------|---------------------|------------------|----------------|
| 12.02   | 2.94 ng | 205284              | Phenanthrene-d10 | 11.52          |
| Hit# of | 5       | Tentative ID        | MW MolForm       | CAS# Qual      |
| 1       |         | n-Hexadecanoic acid | 256 C16H32O2     | 000057-10-3 97 |
| 2       |         | Tridecanoic acid    | 214 C13H26O2     | 000638-53-9 76 |
| 3       |         | Pentadecanoic acid  | 242 C15H30O2     | 001002-84-2 74 |
| 4       |         | Tetradecanoic acid  | 228 C14H28O2     | 000544-63-8 72 |
| 5       |         | n-Decanoic acid     | 172 C10H20O2     | 000334-48-5 53 |



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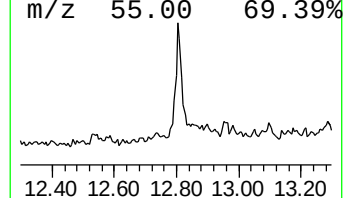
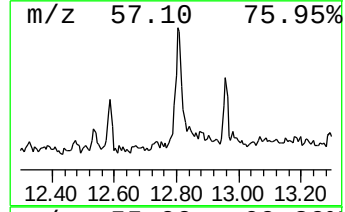
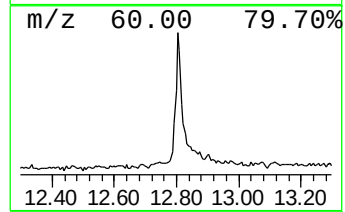
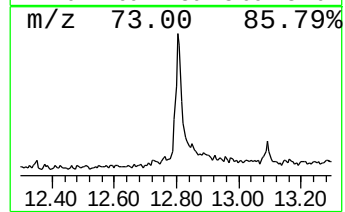
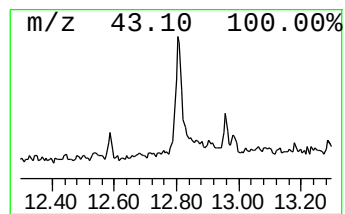
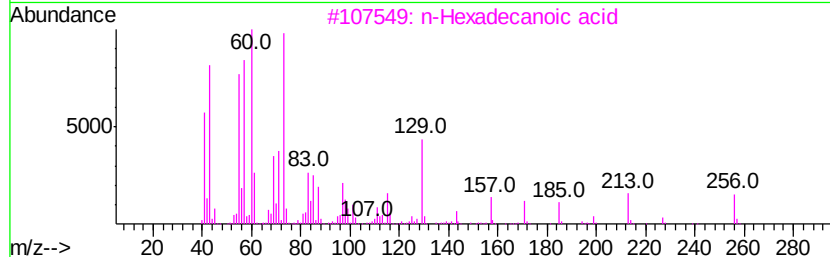
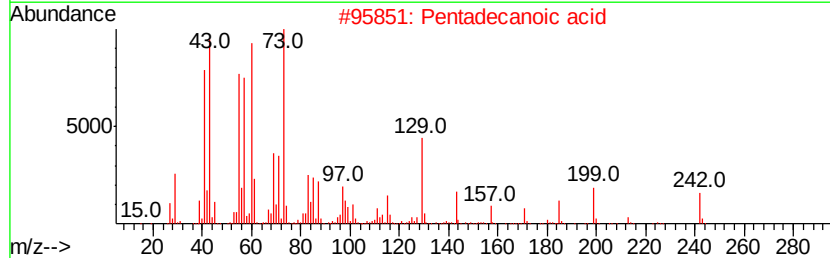
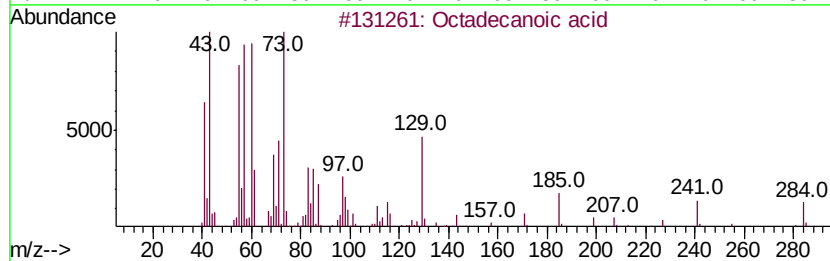
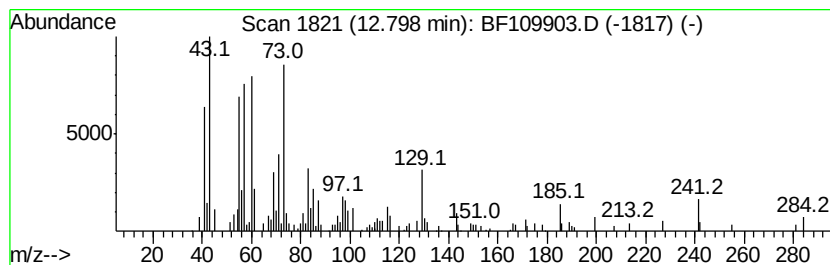
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ClientSampleId :  
AOC3-01A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Octadecanoic acid Concentration Rank 6

| R.T.      | EstConc             | Area   | Relative to ISTD |             | R.T.  |
|-----------|---------------------|--------|------------------|-------------|-------|
| 12.80     | 4.99 ng             | 348140 | Phenanthrene-d10 |             | 11.52 |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual  |
| 1         | Octadecanoic acid   | 284    | C18H36O2         | 000057-11-4 | 99    |
| 2         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 94    |
| 3         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 91    |
| 4         | n-Decanoic acid     | 172    | C10H20O2         | 000334-48-5 | 68    |
| 5         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 64    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\  
Data File : BF109903.D  
Acq On : 16 Oct 2018 14:20  
Operator : JU/SJ  
Sample : J5515-10  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

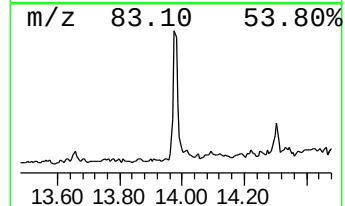
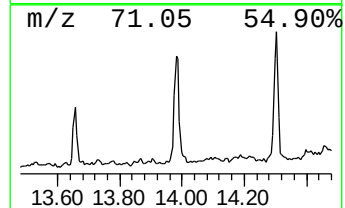
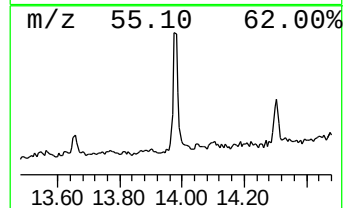
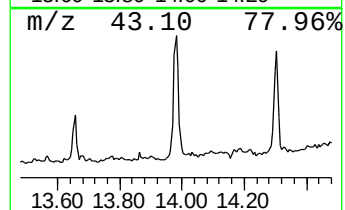
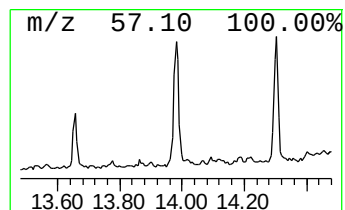
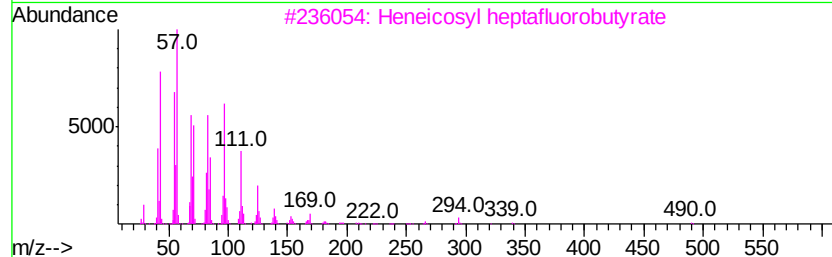
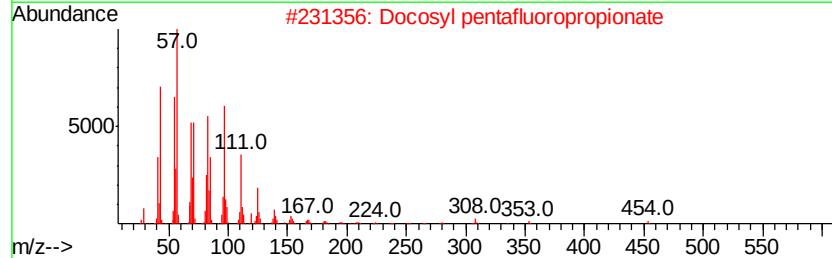
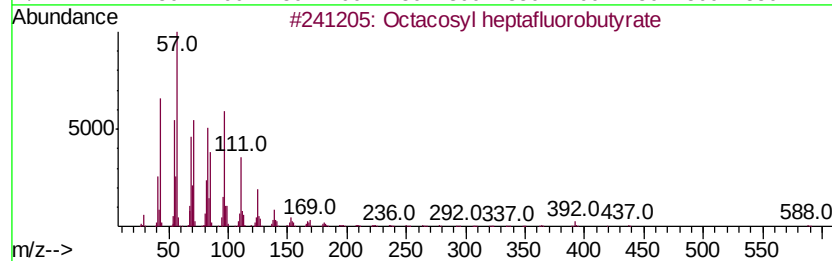
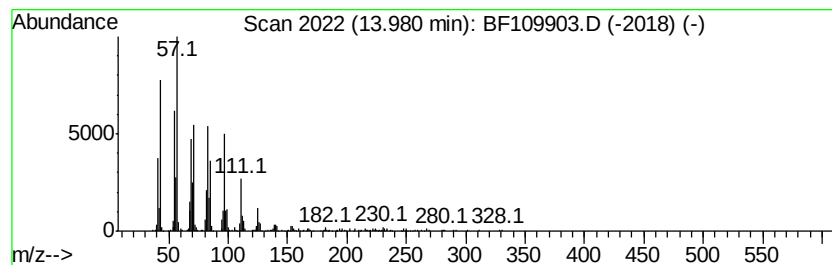
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Octacosyl heptafluorobutyrate Concentration Rank 5

| R.T.      | EstConc                                    | Area   | Relative to ISTD |              | R.T.  |
|-----------|--------------------------------------------|--------|------------------|--------------|-------|
| 13.98     | 5.68 ng                                    | 442579 | Chrysene-d12     |              | 14.16 |
| Hit# of 5 | Tentative ID                               | MW     | MolForm          | CAS#         | Qual  |
| 1         | Octacosyl heptafluorobutyrate              | 606    | C32H57F7O2       | 1000351-83-6 | 91    |
| 2         | Docosyl pentafluoropropionate              | 472    | C25H45F5O2       | 1000351-80-9 | 91    |
| 3         | Heneicosyl heptafluorobutyrate             | 508    | C25H43F7O2       | 1000351-83-8 | 91    |
| 4         | Nonadecyl pentafluoropropionate            | 430    | C22H39F5O2       | 1000351-88-8 | 91    |
| 5         | Pentafluoropropionic acid, octadecyl ester | 416    | C21H37F5O2       | 959261-25-7  | 91    |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\  
Data File : BF109903.D  
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Operator : JU/SJ  
Sample : J5515-10  
Misc :  
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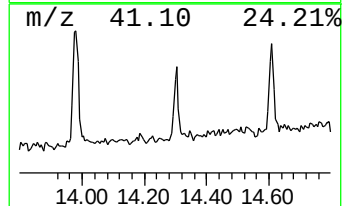
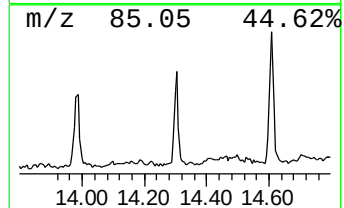
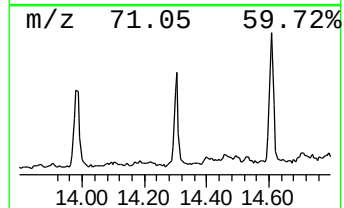
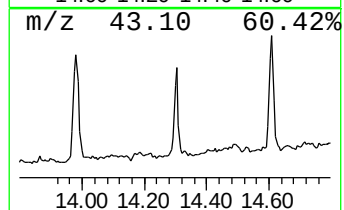
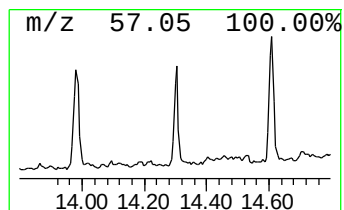
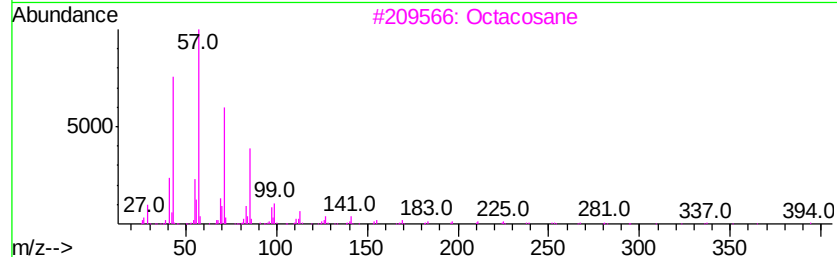
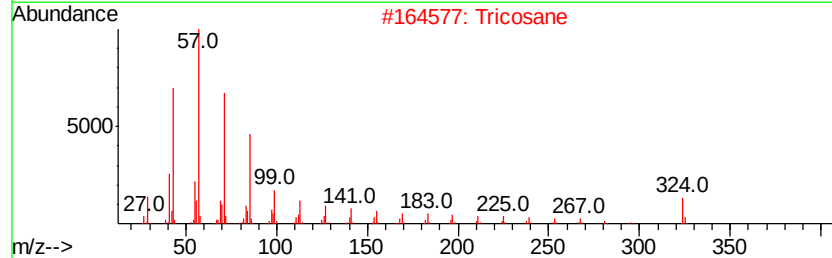
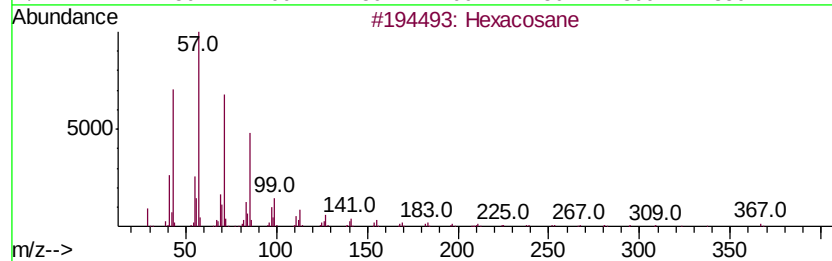
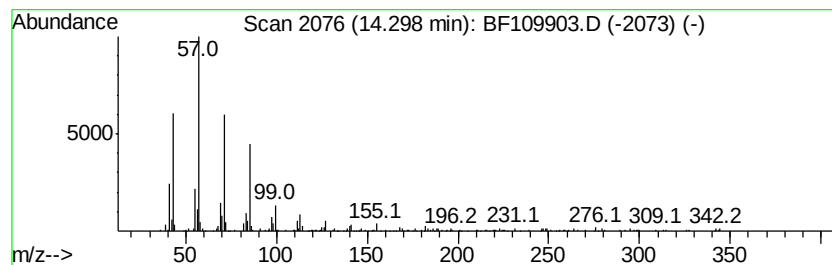
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Hexacosane Concentration Rank 12

| R.T.    | EstConc     | Area         | Relative to ISTD | R.T.      |
|---------|-------------|--------------|------------------|-----------|
| 14.30   | 3.11 ng     | 242376       | Chrysene-d12     | 14.16     |
| Hit# of | 5           | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Hexacosane  | 366 C26H54   | 000630-01-3      | 90        |
| 2       | Tricosane   | 324 C23H48   | 000638-67-5      | 90        |
| 3       | Octacosane  | 394 C28H58   | 000630-02-4      | 90        |
| 4       | Heneicosane | 296 C21H44   | 000629-94-7      | 90        |
| 5       | Octadecane  | 254 C18H38   | 000593-45-3      | 87        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\  
Data File : BF109903.D  
Acq On : 16 Oct 2018 14:20  
Operator : JU/SJ  
Sample : J5515-10  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

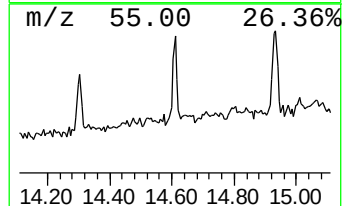
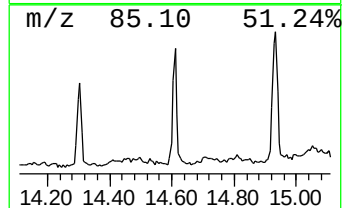
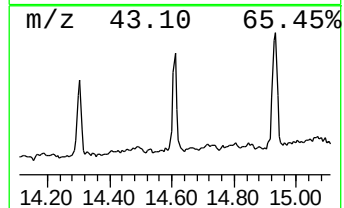
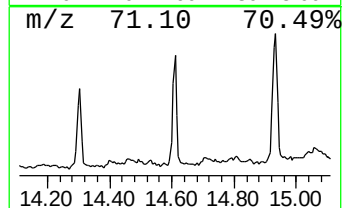
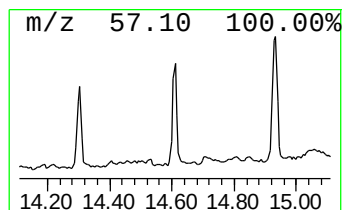
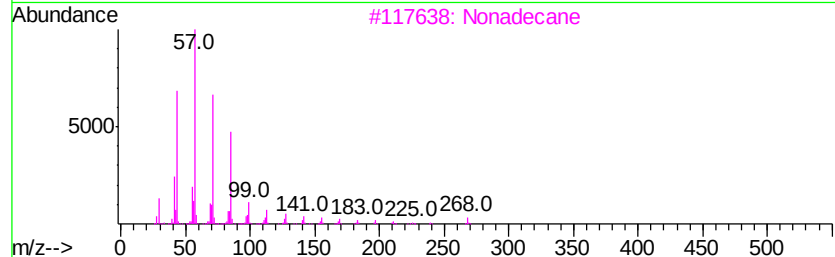
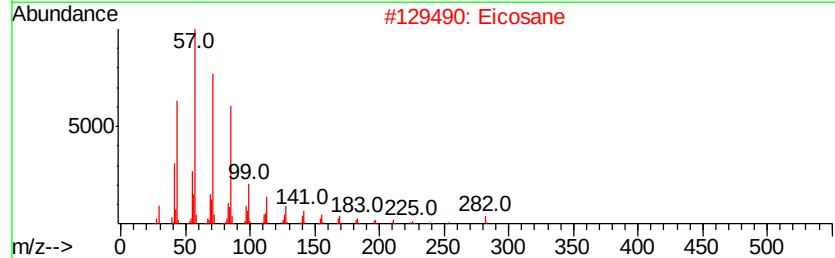
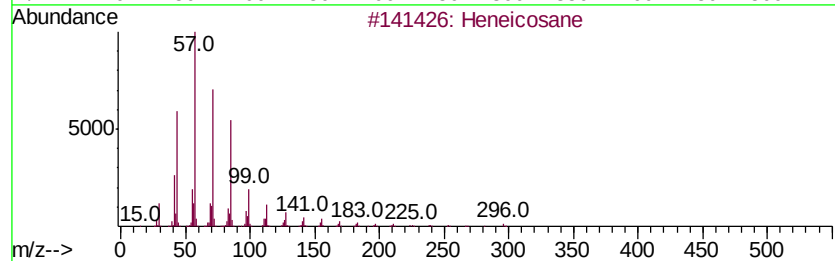
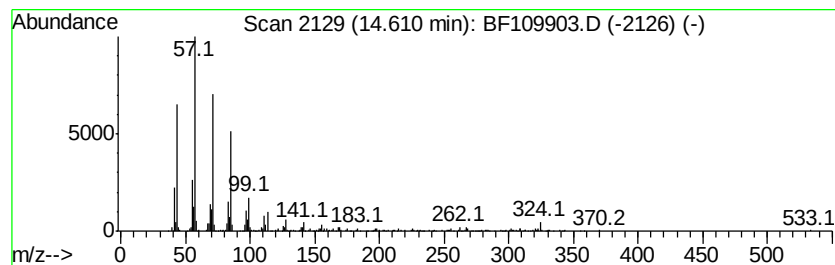
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Heneicosane Concentration Rank 11

| R.T.    | EstConc     | Area         | Relative to ISTD | R.T.           |
|---------|-------------|--------------|------------------|----------------|
| 14.61   | 3.86 ng     | 300646       | Chrysene-d12     | 14.16          |
| Hit# of | 5           | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Heneicosane |              | 296 C21H44       | 000629-94-7 97 |
| 2       | Eicosane    |              | 282 C20H42       | 000112-95-8 95 |
| 3       | Nonadecane  |              | 268 C19H40       | 000629-92-5 91 |
| 4       | Hexacosane  |              | 366 C26H54       | 000630-01-3 90 |
| 5       | Heptadecane |              | 240 C17H36       | 000629-78-7 87 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\  
Data File : BF109903.D  
Acq On : 16 Oct 2018 14:20  
Operator : JU/SJ  
Sample : J5515-10  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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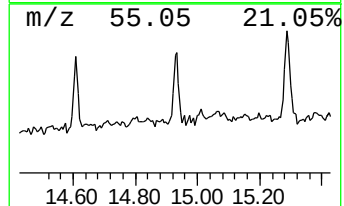
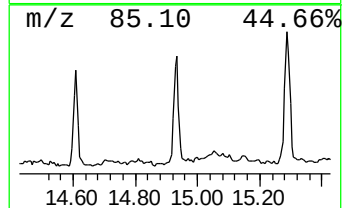
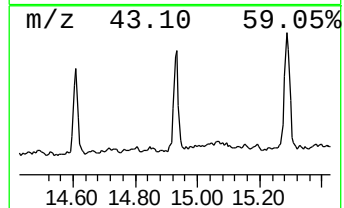
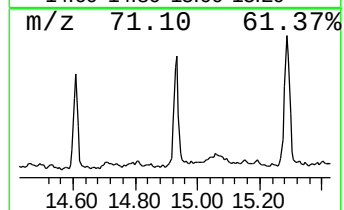
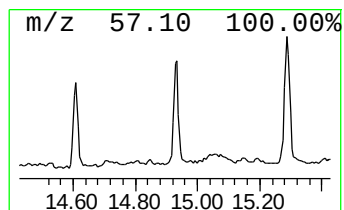
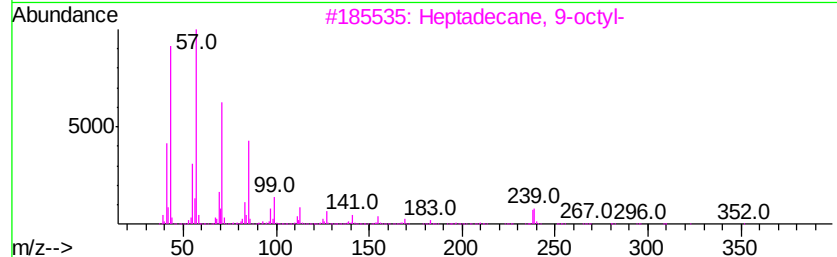
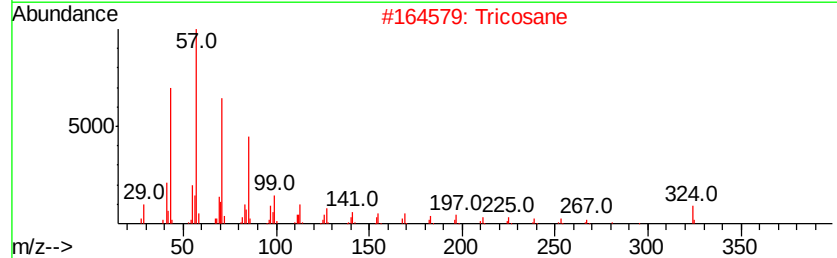
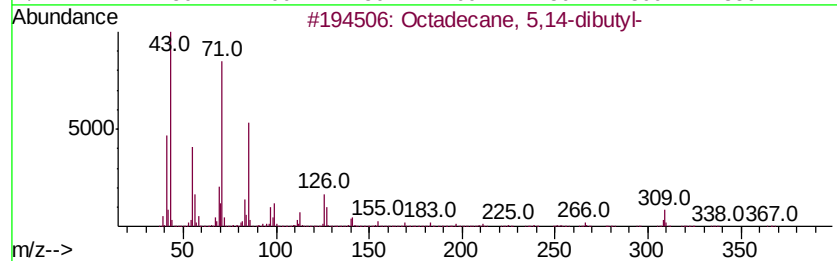
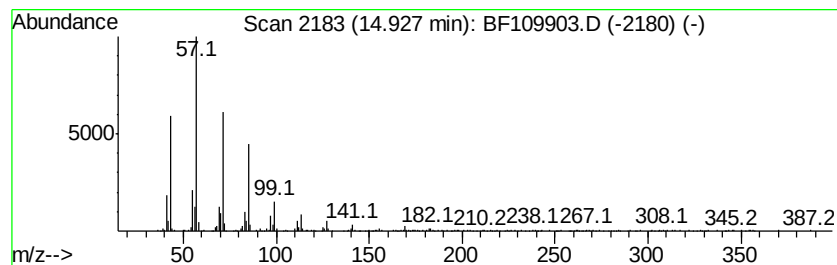
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Octadecane, 5,14-dibutyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.93 | 3.91 ng | 413273 | Perylene-d12     | 15.69 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Octadecane, 5,14-dibutyl- | 366 | C26H54  | 055282-13-8 | 93   |
| 2    |    | Tricosane                 | 324 | C23H48  | 000638-67-5 | 93   |
| 3    |    | Heptadecane, 9-octyl-     | 352 | C25H52  | 007225-64-1 | 93   |
| 4    |    | Hexacosane                | 366 | C26H54  | 000630-01-3 | 91   |
| 5    |    | Tetracosane               | 338 | C24H50  | 000646-31-1 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\  
Data File : BF109903.D  
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Operator : JU/SJ  
Sample : J5515-10  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
AOC3-01A

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

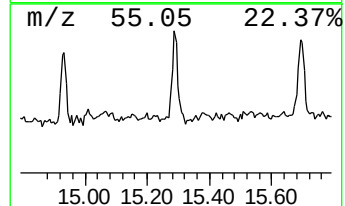
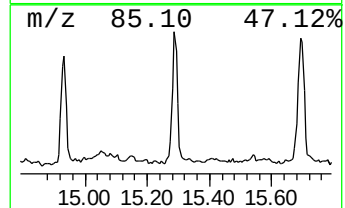
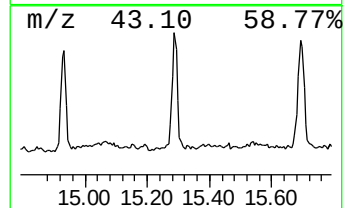
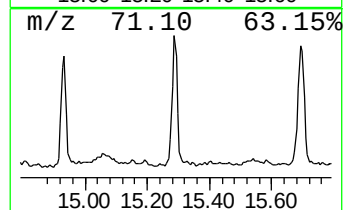
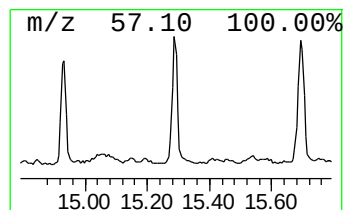
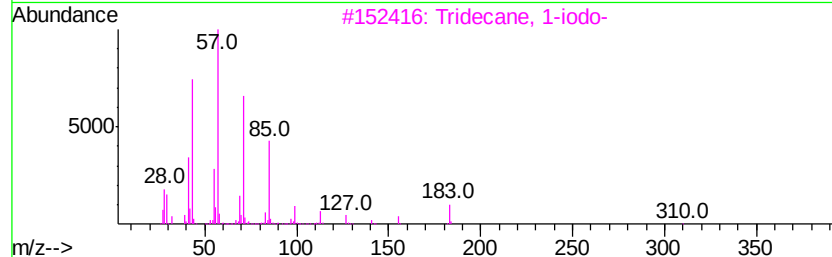
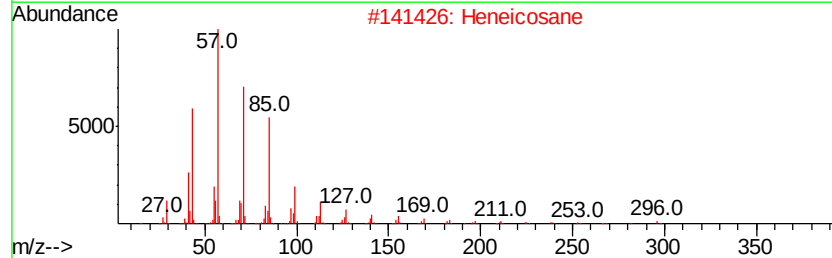
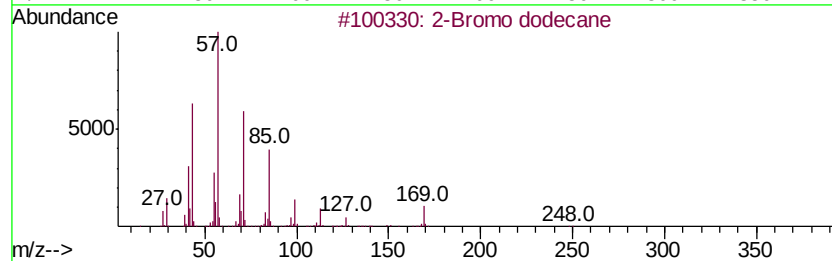
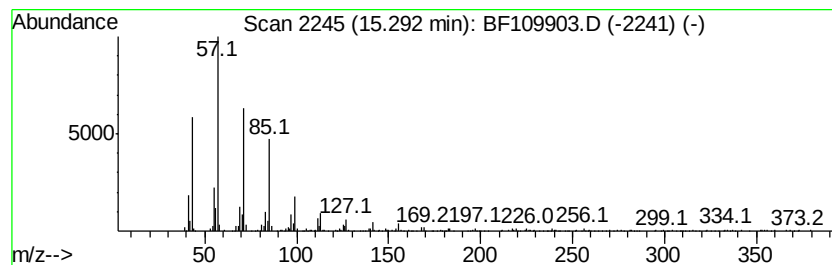
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 2-Bromo dodecane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.29 | 4.55 ng | 481809 | Perylene-d12     | 15.69 |

| Hit# | of | Tentative ID          | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Bromo dodecane      | 248 | C12H25Br | 013187-99-0 | 93   |
| 2    |    | Heneicosane           | 296 | C21H44   | 000629-94-7 | 91   |
| 3    |    | Tridecane, 1-iodo-    | 310 | C13H27I  | 035599-77-0 | 91   |
| 4    |    | Octacosane            | 394 | C28H58   | 000630-02-4 | 91   |
| 5    |    | Heptadecane, 9-octyl- | 352 | C25H52   | 007225-64-1 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101618\  
Data File : BF109903.D  
Acq On : 16 Oct 2018 14:20  
Operator : JU/SJ  
Sample : J5515-10  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
AOC3-01A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

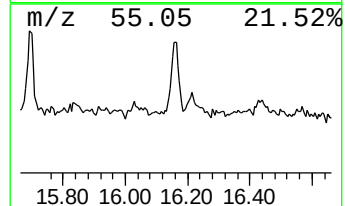
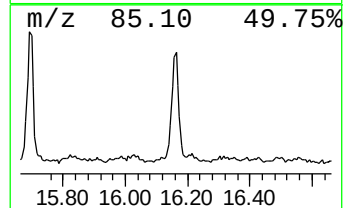
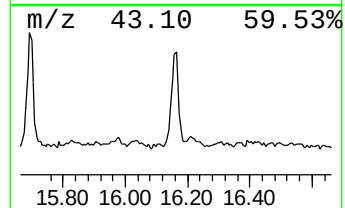
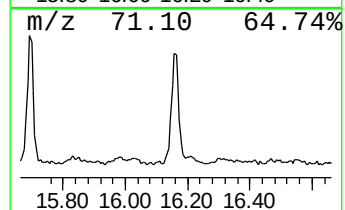
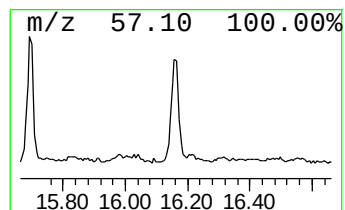
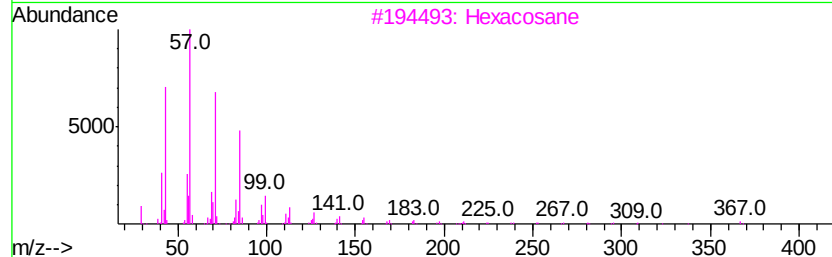
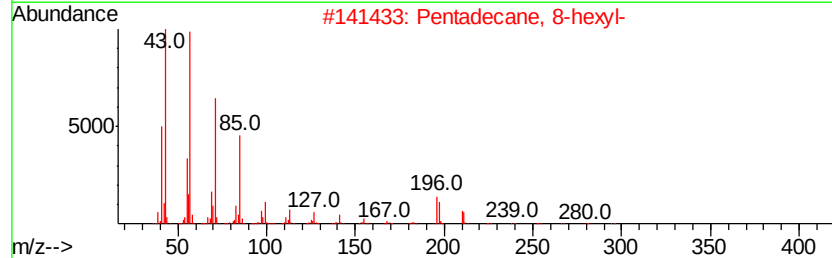
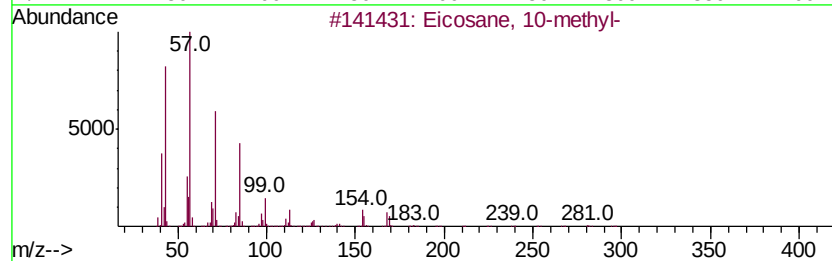
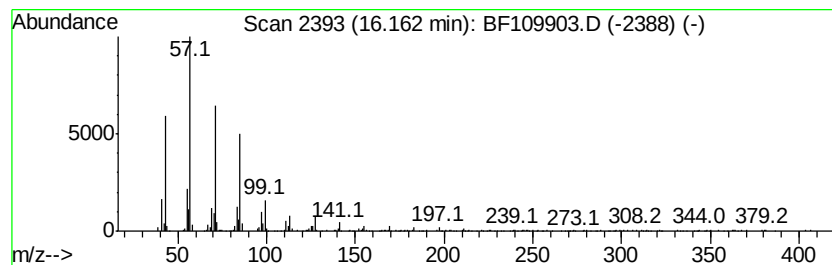
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Eicosane, 10-methyl- Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.16 | 4.65 ng | 492012 | Perylene-d12     | 15.69 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane, 10-methyl-  | 296 | C21H44  | 054833-23-7 | 93   |
| 2    |    | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 93   |
| 3    |    | Hexacosane            | 366 | C26H54  | 000630-01-3 | 93   |
| 4    |    | Tridecane, 6-propyl-  | 226 | C16H34  | 055045-10-8 | 90   |
| 5    |    | Tetracosane           | 338 | C24H50  | 000646-31-1 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF101618\  
Data File : BF109903.D  
Acq On : 16 Oct 2018 14:20  
Operator : JU/SJ  
Sample : J5515-10  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
AOC3-01A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF101518.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Butane, 2-methoxy... | 2.33  | 12.4    | ng    | 912034   | 1                     | 6.98  | 1472890 | 20.0 |
| 2-Pentanone, 4-hy... | 5.22  | 2.7     | ng    | 195942   | 1                     | 6.98  | 1472890 | 20.0 |
| unknown6.75          | 6.75  | 67.4    | ng    | 4962650  | 1                     | 6.98  | 1472890 | 20.0 |
| n-Hexadecanoic acid  | 12.02 | 2.9     | ng    | 205284   | 4                     | 11.52 | 1395540 | 20.0 |
| Octadecanoic acid    | 12.80 | 5.0     | ng    | 348140   | 4                     | 11.52 | 1395540 | 20.0 |
| Octacosyl heptafl... | 13.98 | 5.7     | ng    | 442579   | 5                     | 14.16 | 1558610 | 20.0 |
| Hexacosane           | 14.30 | 3.1     | ng    | 242376   | 5                     | 14.16 | 1558610 | 20.0 |
| Heneicosane          | 14.61 | 3.9     | ng    | 300646   | 5                     | 14.16 | 1558610 | 20.0 |
| Octadecane, 5,14-... | 14.93 | 3.9     | ng    | 413273   | 6                     | 15.69 | 2115620 | 20.0 |
| 2-Bromo dodecane     | 15.29 | 4.5     | ng    | 481809   | 6                     | 15.69 | 2115620 | 20.0 |
| Eicosane, 10-methyl- | 16.16 | 4.7     | ng    | 492012   | 6                     | 15.69 | 2115620 | 20.0 |