

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF040820\  
 Data File : BF119859.D  
 Acq On : 9 Apr 2020 7:38  
 Operator : MA/SJ  
 Sample : L2151-09  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MH-109

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-BF040820.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.963       | 485           | 489         | 496          | rVB      | 105269         | 119962        | 1.33%           | 0.241%        |
| 2         | 5.375       | 554           | 559         | 562          | rBV      | 3169731        | 2821227       | 31.37%          | 5.664%        |
| 3         | 6.392       | 727           | 732         | 735          | rBV      | 1899897        | 1653131       | 18.38%          | 3.319%        |
| 4         | 6.763       | 791           | 795         | 799          | rBV      | 1256017        | 1068640       | 11.88%          | 2.146%        |
| 5         | 6.922       | 817           | 822         | 826          | rBV      | 4690813        | 4879762       | 54.25%          | 9.798%        |
| 6         | 7.334       | 885           | 892         | 895          | rBV      | 4129448        | 4678371       | 52.01%          | 9.393%        |
| 7         | 8.051       | 1009          | 1014        | 1017         | rBV      | 1579240        | 1411486       | 15.69%          | 2.834%        |
| 8         | 9.133       | 1192          | 1198        | 1201         | rBV      | 7437305        | 8060585       | 89.62%          | 16.184%       |
| 9         | 9.522       | 1260          | 1264        | 1267         | rBV      | 536895         | 428085        | 4.76%           | 0.860%        |
| 10        | 9.810       | 1306          | 1313        | 1316         | rBV      | 1865328        | 1752091       | 19.48%          | 3.518%        |
| 11        | 10.604      | 1441          | 1448        | 1451         | rBV      | 5182839        | 5688853       | 63.25%          | 11.422%       |
| 12        | 11.292      | 1561          | 1565        | 1569         | rBV      | 2178750        | 1871271       | 20.81%          | 3.757%        |
| 13        | 11.827      | 1652          | 1656        | 1665         | rBV      | 796329         | 728862        | 8.10%           | 1.463%        |
| 14        | 12.610      | 1786          | 1789        | 1801         | rBV      | 244134         | 300338        | 3.34%           | 0.603%        |
| 15        | 12.892      | 1831          | 1837        | 1840         | rBV      | 7761535        | 8994298       | 100.00%         | 18.059%       |
| 16        | 13.363      | 1914          | 1917        | 1921         | rVB      | 132445         | 121193        | 1.35%           | 0.243%        |
| 17        | 13.792      | 1986          | 1990        | 2003         | rBV      | 479935         | 673534        | 7.49%           | 1.352%        |
| 18        | 13.933      | 2010          | 2014        | 2018         | rBV      | 1727718        | 1497508       | 16.65%          | 3.007%        |
| 19        | 14.104      | 2040          | 2043        | 2046         | rBV      | 138582         | 130012        | 1.45%           | 0.261%        |
| 20        | 14.410      | 2091          | 2095        | 2098         | rBV      | 193551         | 180065        | 2.00%           | 0.362%        |
| 21        | 14.521      | 2111          | 2114        | 2119         | rVB      | 126067         | 110981        | 1.23%           | 0.223%        |
| 22        | 14.710      | 2143          | 2146        | 2150         | rBV      | 201311         | 213834        | 2.38%           | 0.429%        |
| 23        | 15.039      | 2198          | 2202        | 2208         | rBV      | 178476         | 194157        | 2.16%           | 0.390%        |
| 24        | 15.368      | 2253          | 2258        | 2262         | rBV      | 993559         | 1197283       | 13.31%          | 2.404%        |
| 25        | 15.410      | 2262          | 2265        | 2273         | rVB      | 153744         | 185054        | 2.06%           | 0.372%        |
| 26        | 15.833      | 2331          | 2337        | 2343         | rVB      | 109096         | 161162        | 1.79%           | 0.324%        |
| 27        | 16.315      | 2414          | 2419        | 2431         | rVB3     | 74418          | 154398        | 1.72%           | 0.310%        |
| 28        | 17.580      | 2616          | 2634        | 2636         | rBV3     | 141195         | 529765        | 5.89%           | 1.064%        |

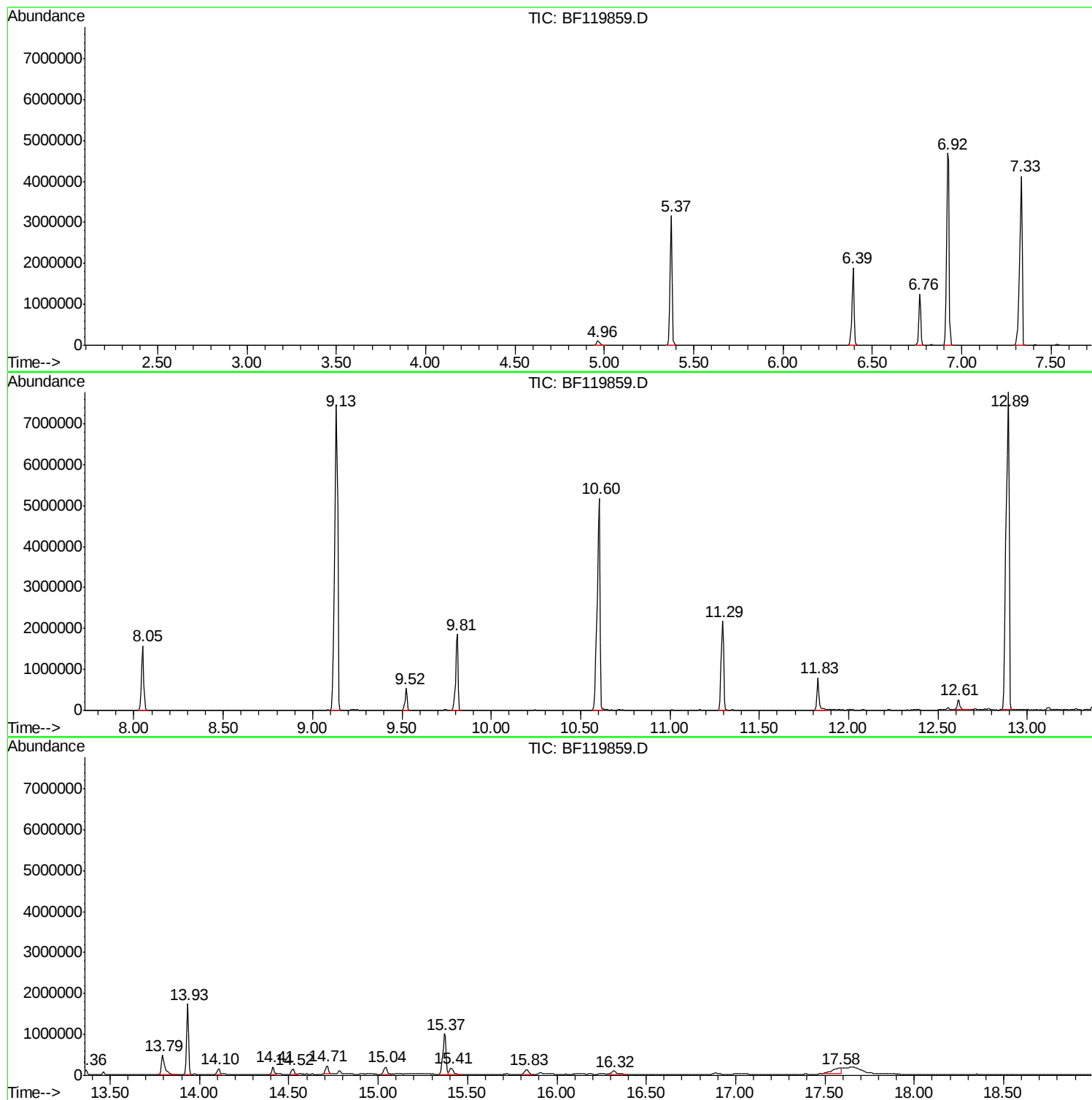
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ClientSampleId :  
MH-109

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.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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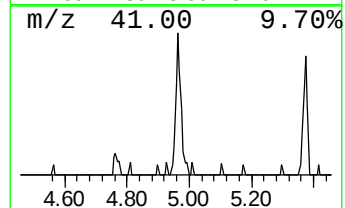
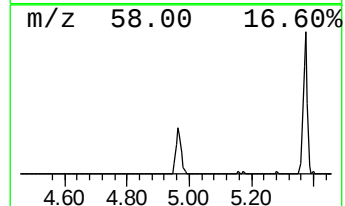
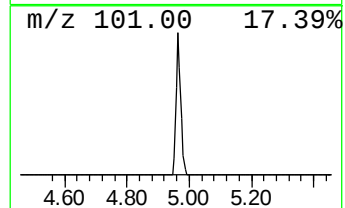
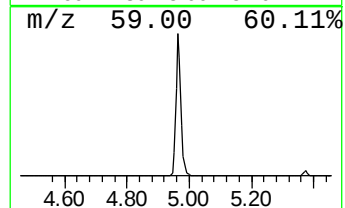
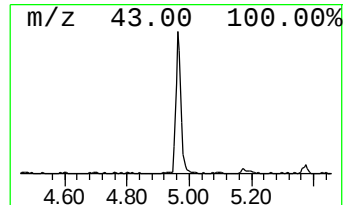
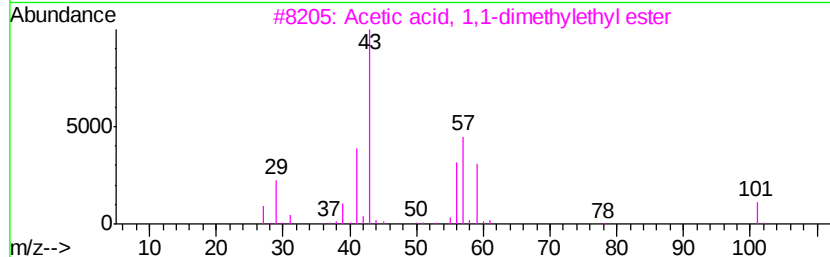
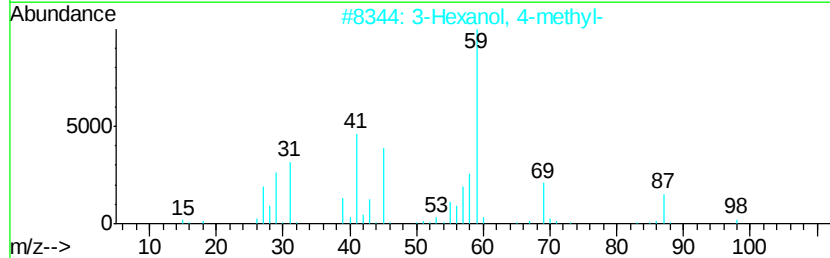
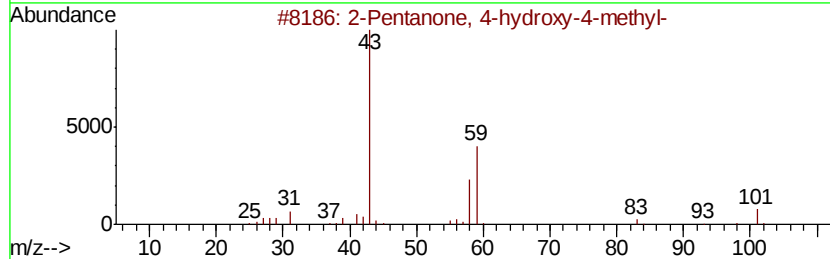
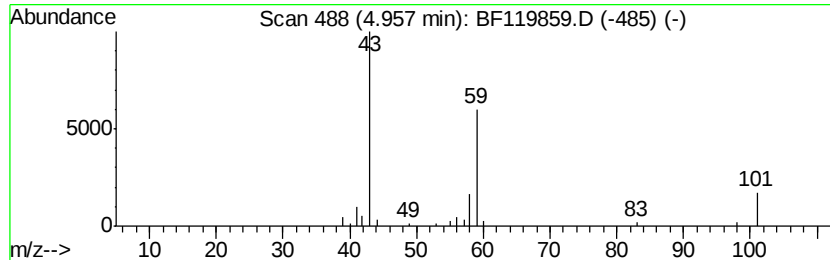
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.96 | 2.25 ng | 119962 | 1,4-Dichlorobenzene-d4 | 6.76 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    |   | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2 | 33   |
| 3    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 28   |
| 4    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 5    |    |   | 2,3-Butanedione, monooxime          | 101 | C4H7NO2  | 000057-71-6 | 9    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

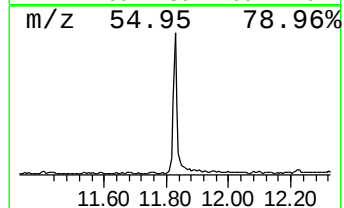
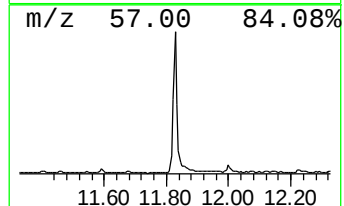
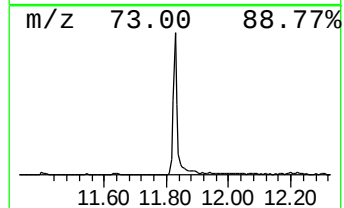
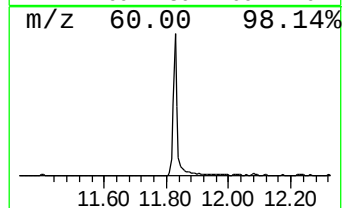
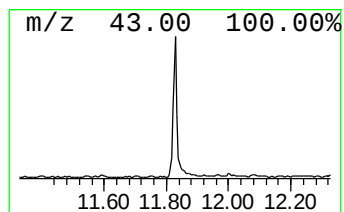
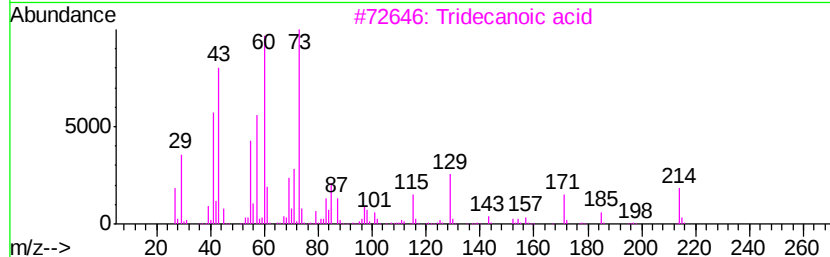
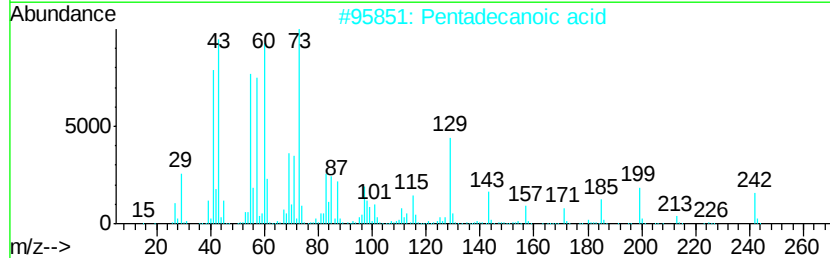
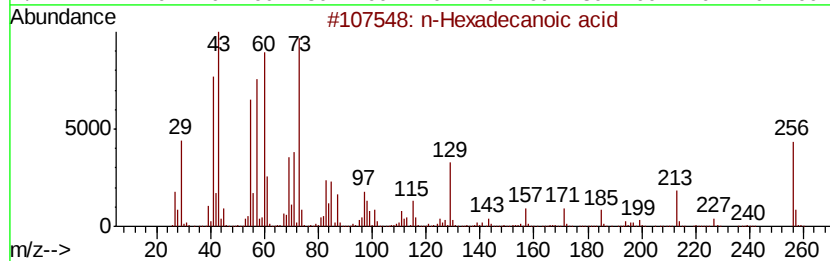
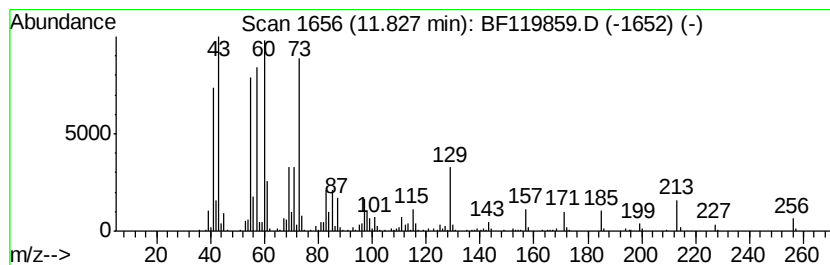
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.83 | 7.79 ng | 728862 | Phenanthrene-d10 | 11.29 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 95   |
| 2    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 87   |
| 3    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 83   |
| 4    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 78   |
| 5    |    | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 64   |



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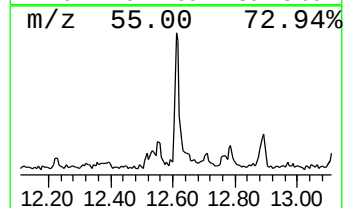
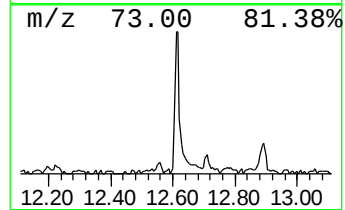
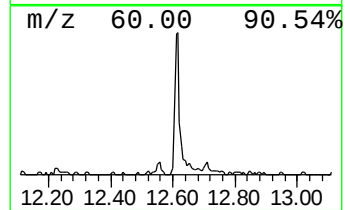
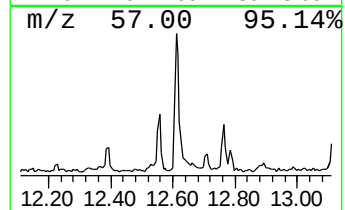
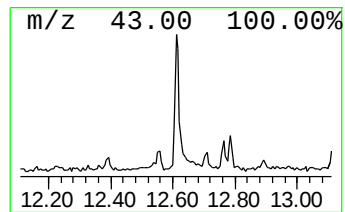
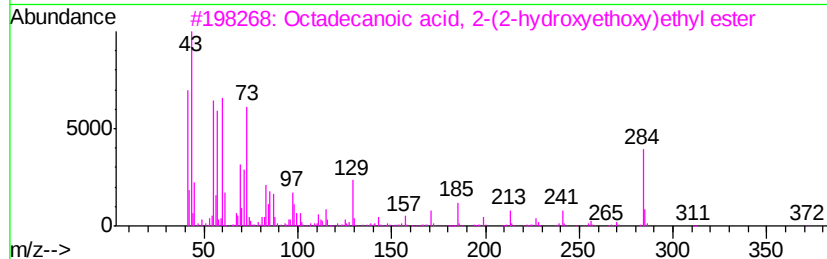
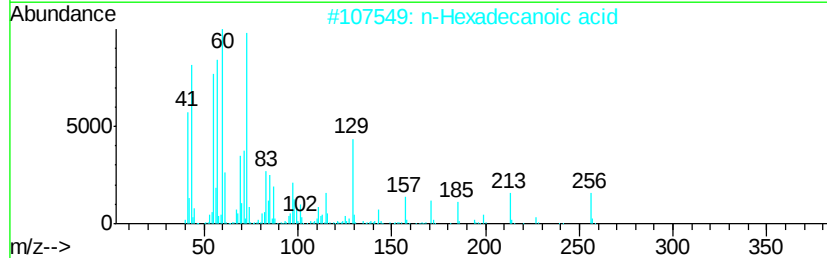
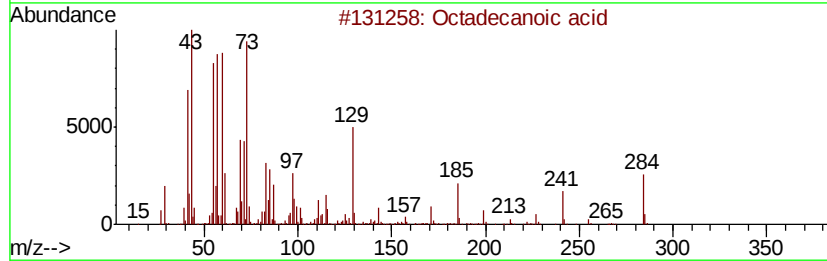
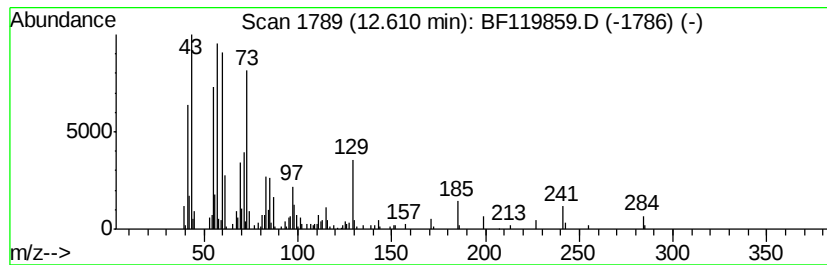
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Octadecanoic acid Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.61 | 3.21 ng | 300338 | Phenanthrene-d10 | 11.29 |

| Hit# of | 5                                    | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|---------|--------------------------------------|--------------|-----|----------|-------------|------|
| 1       | Octadecanoic acid                    |              | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2       | n-Hexadecanoic acid                  |              | 256 | C16H32O2 | 000057-10-3 | 87   |
| 3       | Octadecanoic acid, 2-(2-hydroxyve... |              | 372 | C22H44O4 | 000106-11-6 | 86   |
| 4       | n-Decanoic acid                      |              | 172 | C10H20O2 | 000334-48-5 | 62   |
| 5       | Pentadecanoic acid                   |              | 242 | C15H30O2 | 001002-84-2 | 53   |



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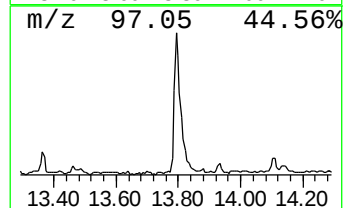
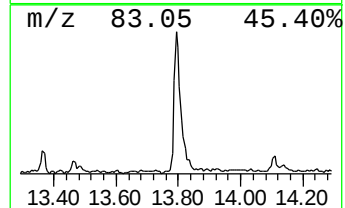
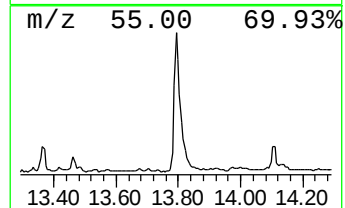
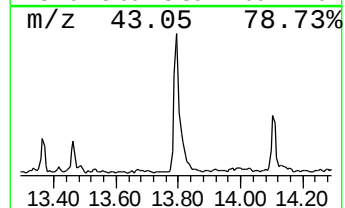
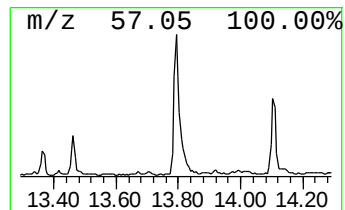
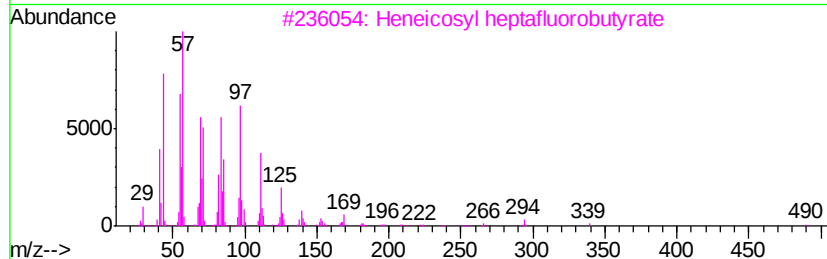
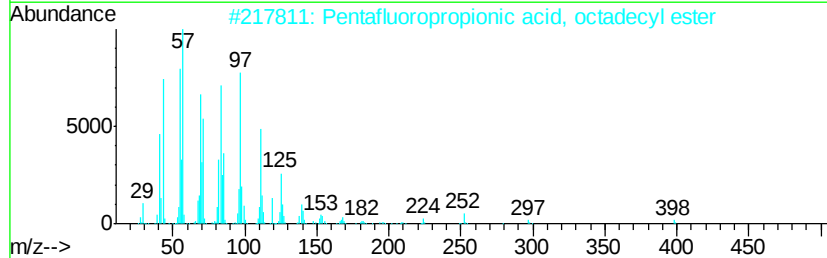
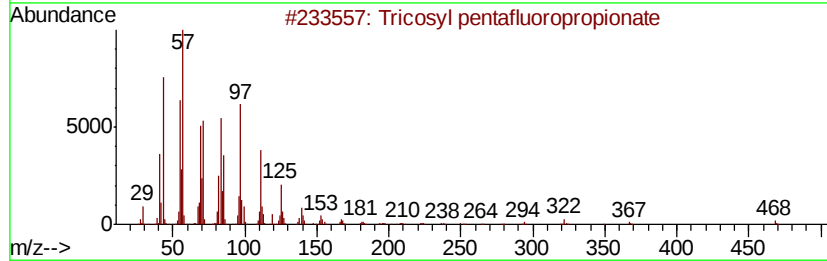
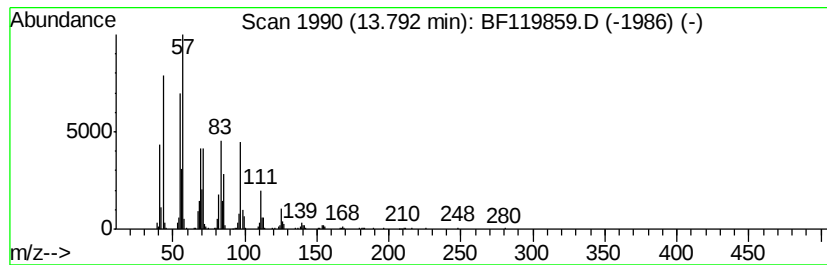
TIC Library : C:\DATABASE\NIST11.L  
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Peak Number 5 Tricosyl pentafluoropropionate Concentration Rank 2

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.79 | 9.00 ng | 673534 | Chrysene-d12     | 13.93 |

| Hit# | of | Tentative ID                               | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Tricosyl pentafluoropropionate             | 486 | C26H47F5O2 | 1000351-81-0 | 91   |
| 2    |    | Pentafluoropropionic acid, octadecyl ester | 416 | C21H37F5O2 | 959261-25-7  | 91   |
| 3    |    | Heneicosyl heptafluorobutyrate             | 508 | C25H43F7O2 | 1000351-83-8 | 91   |
| 4    |    | Heptafluorobutyric acid, hexadecyl ester   | 438 | C20H33F7O2 | 006385-15-5  | 91   |
| 5    |    | Heptadecyl trifluoroacetate                | 352 | C19H35F3O2 | 1000351-87-0 | 91   |



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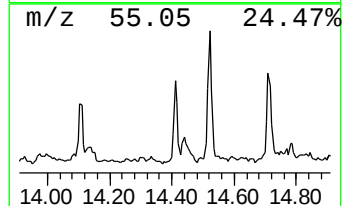
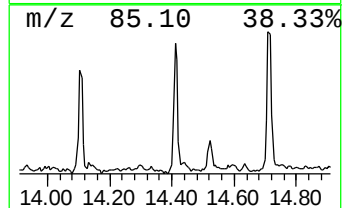
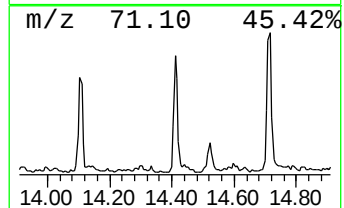
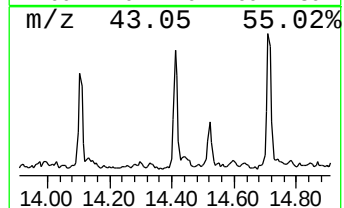
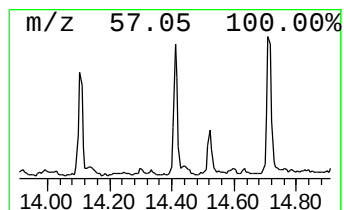
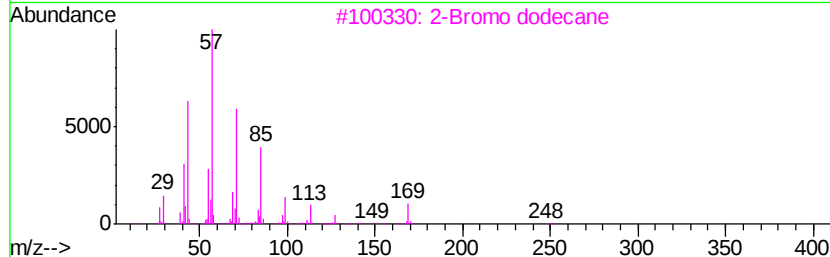
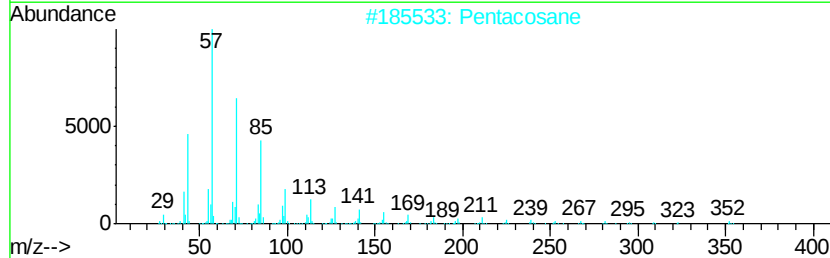
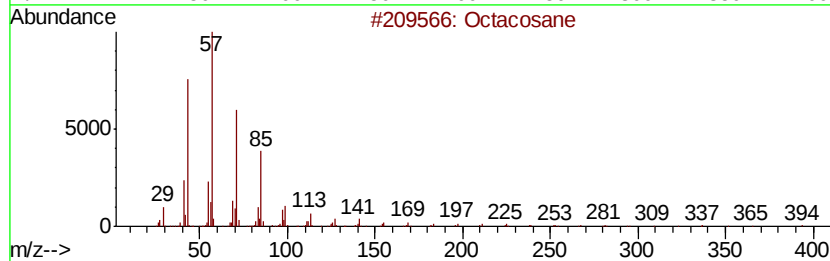
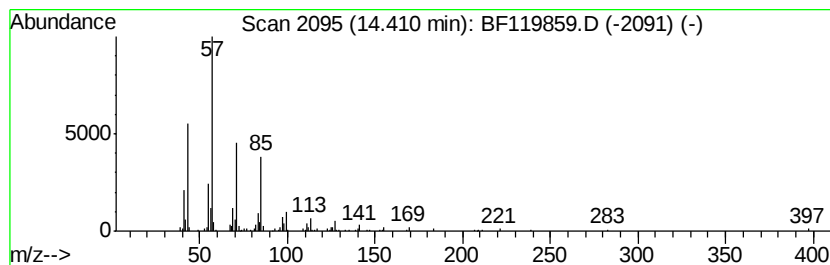
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Octacosane Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.41 | 2.40 ng | 180065 | Chrysene-d12     | 13.93 |

| Hit# of | 5                  | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|---------|--------------------|--------------|-----|----------|-------------|------|
| 1       | Octacosane         |              | 394 | C28H58   | 000630-02-4 | 91   |
| 2       | Pentacosane        |              | 352 | C25H52   | 000629-99-2 | 90   |
| 3       | 2-Bromo dodecane   |              | 248 | C12H25Br | 013187-99-0 | 86   |
| 4       | Tridecane, 1-iodo- |              | 310 | C13H27I  | 035599-77-0 | 86   |
| 5       | Heneicosane        |              | 296 | C21H44   | 000629-94-7 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\  
Data File : BF119859.D  
Acq On : 9 Apr 2020 7:38  
Operator : MA/SJ  
Sample : L2151-09  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MH-109

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

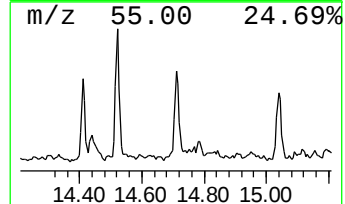
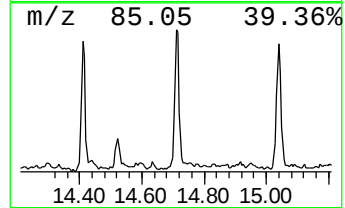
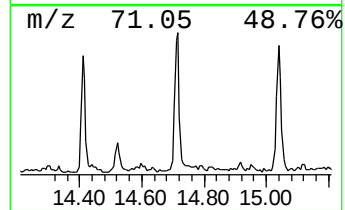
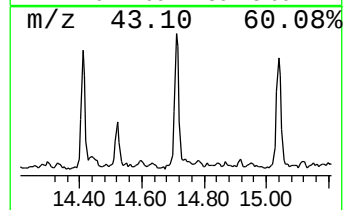
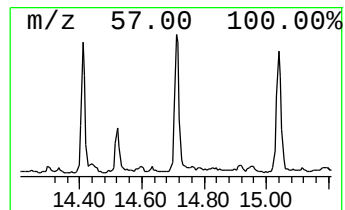
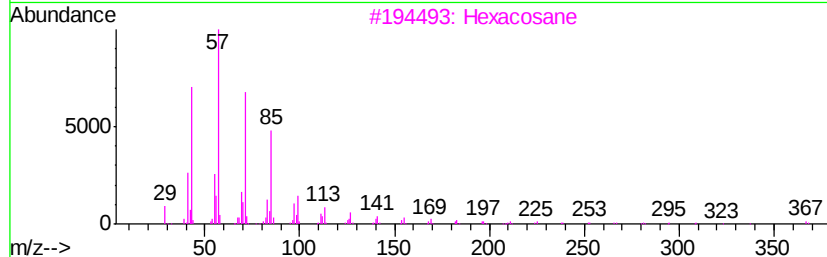
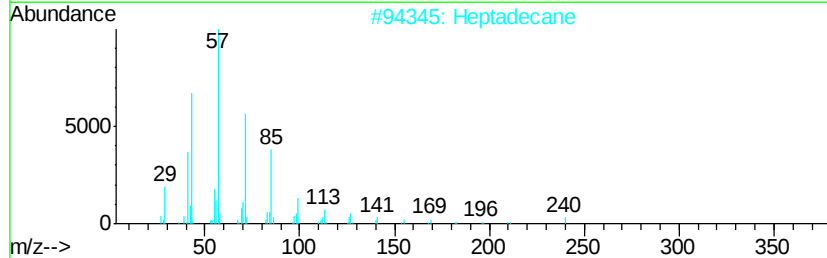
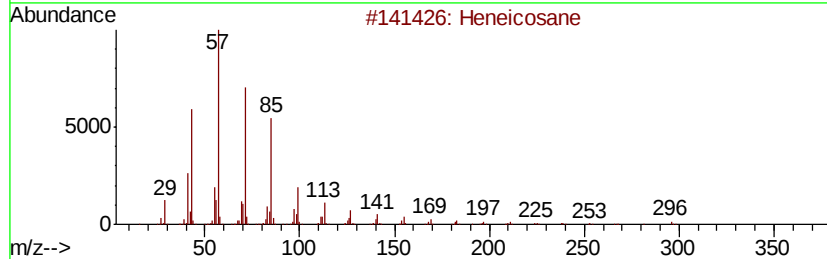
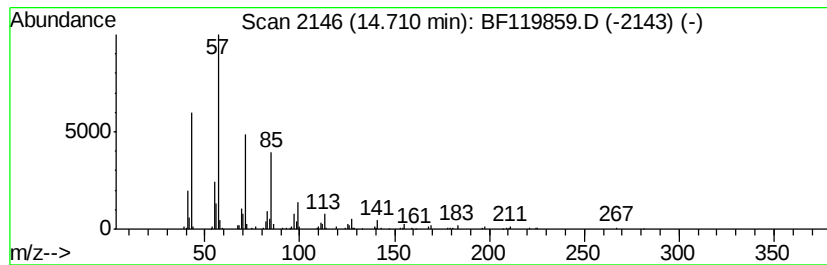
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Heneicosane Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.71 | 3.57 ng | 213834 | Perylene-d12     | 15.37 |

| Hit# of | 5                  | Tentative ID | MW  | MolForm | CAS#         | Qual |
|---------|--------------------|--------------|-----|---------|--------------|------|
| 1       | Heneicosane        |              | 296 | C21H44  | 000629-94-7  | 91   |
| 2       | Heptadecane        |              | 240 | C17H36  | 000629-78-7  | 91   |
| 3       | Hexacosane         |              | 366 | C26H54  | 000630-01-3  | 90   |
| 4       | 2-methyloctacosane |              | 408 | C29H60  | 1000376-72-8 | 87   |
| 5       | Octadecane         |              | 254 | C18H38  | 000593-45-3  | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\  
Data File : BF119859.D  
Acq On : 9 Apr 2020 7:38  
Operator : MA/SJ  
Sample : L2151-09  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MH-109

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

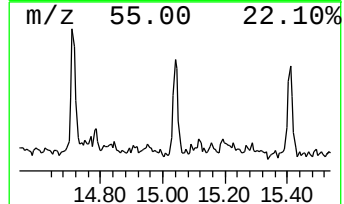
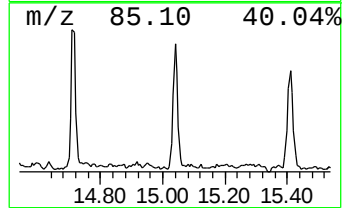
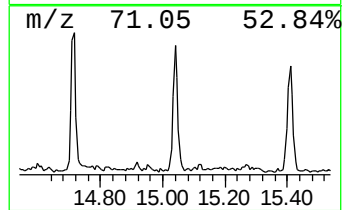
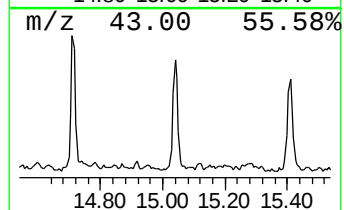
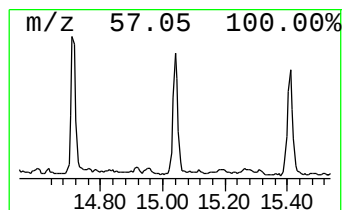
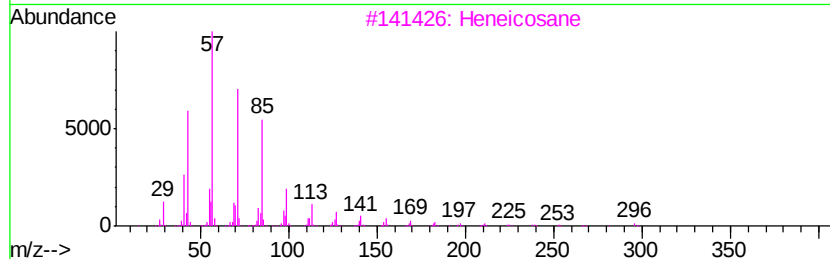
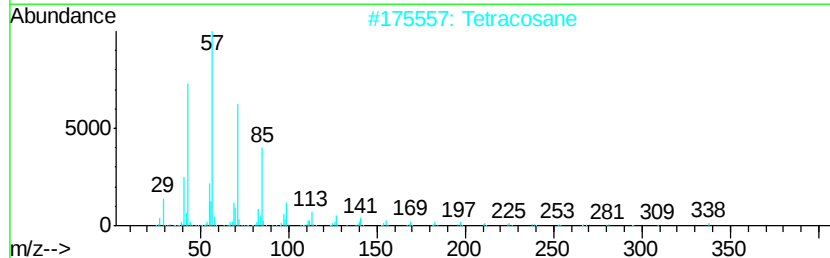
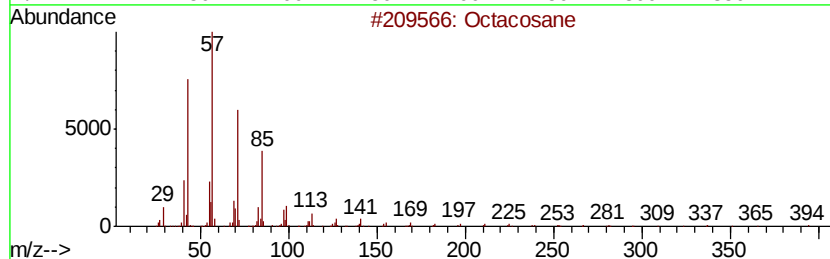
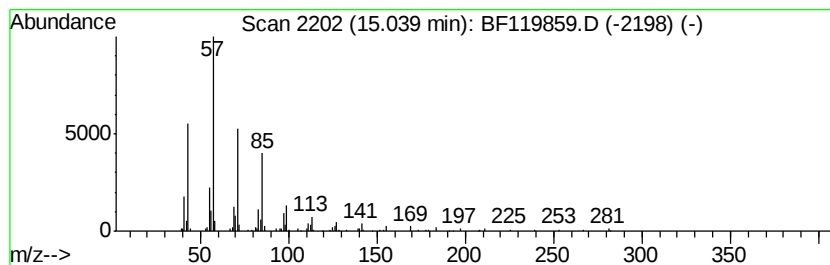
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Tetracosane Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.04 | 3.24 ng | 194157 | Perylene-d12     | 15.37 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Octacosane   | 394 | C28H58  | 000630-02-4 | 91   |
| 2    |    | Tetracosane  | 338 | C24H50  | 000646-31-1 | 91   |
| 3    |    | Heneicosane  | 296 | C21H44  | 000629-94-7 | 90   |
| 4    |    | Heptacosane  | 380 | C27H56  | 000593-49-7 | 90   |
| 5    |    | Hexacosane   | 366 | C26H54  | 000630-01-3 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\  
Data File : BF119859.D  
Acq On : 9 Apr 2020 7:38  
Operator : MA/SJ  
Sample : L2151-09  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MH-109

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

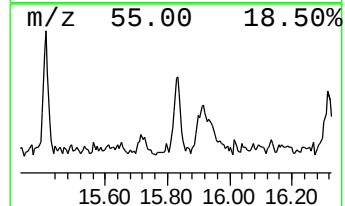
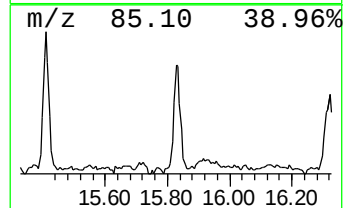
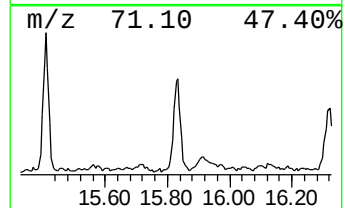
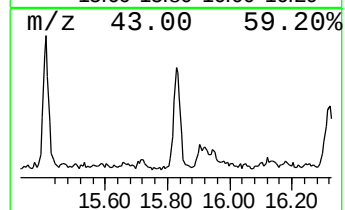
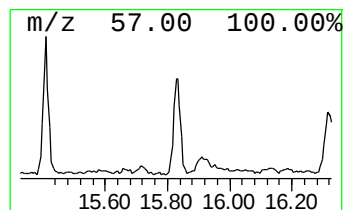
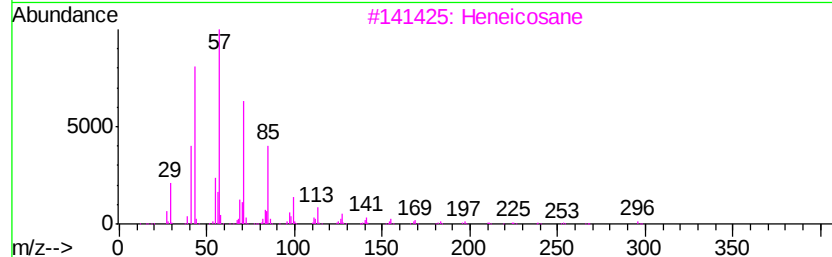
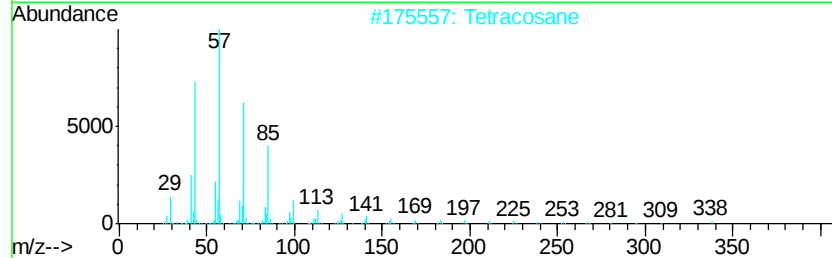
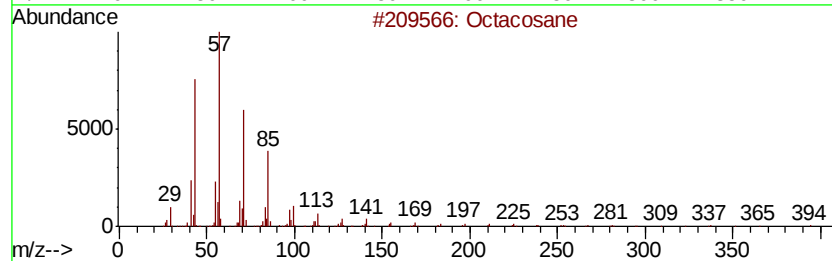
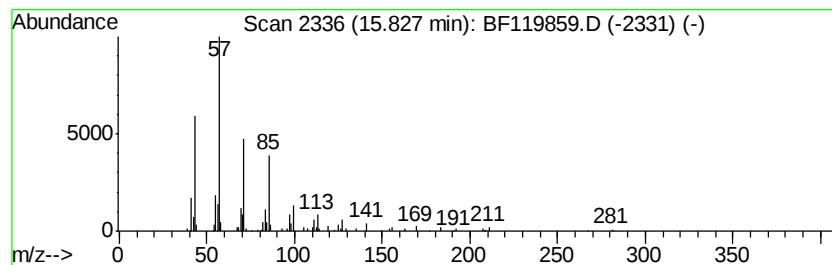
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 2-methyloctacosane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.83 | 2.69 ng | 161162 | Perylene-d12     | 15.37 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------|-----|---------|--------------|------|
| 1    | 5  | Octacosane         | 394 | C28H58  | 000630-02-4  | 91   |
| 2    |    | Tetracosane        | 338 | C24H50  | 000646-31-1  | 91   |
| 3    |    | Heneicosane        | 296 | C21H44  | 000629-94-7  | 87   |
| 4    |    | 2-methyloctacosane | 408 | C29H60  | 1000376-72-8 | 87   |
| 5    |    | Nonadecane         | 268 | C19H40  | 000629-92-5  | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\  
Data File : BF119859.D  
Acq On : 9 Apr 2020 7:38  
Operator : MA/SJ  
Sample : L2151-09  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MH-109

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

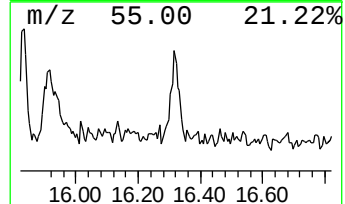
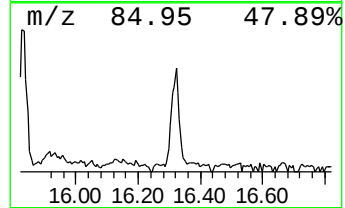
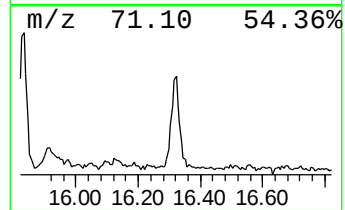
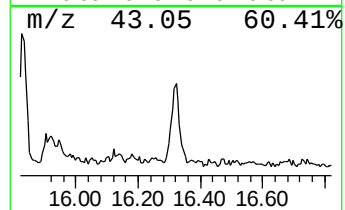
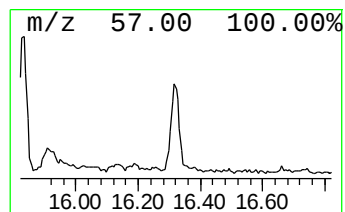
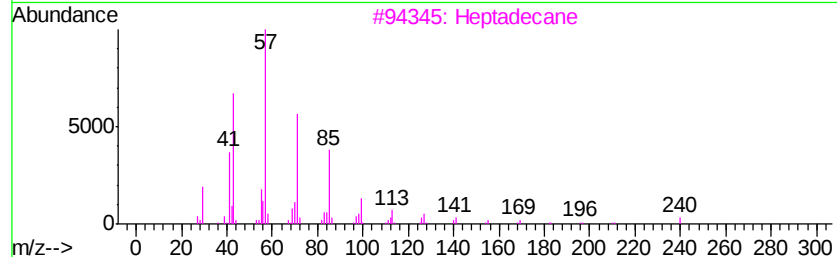
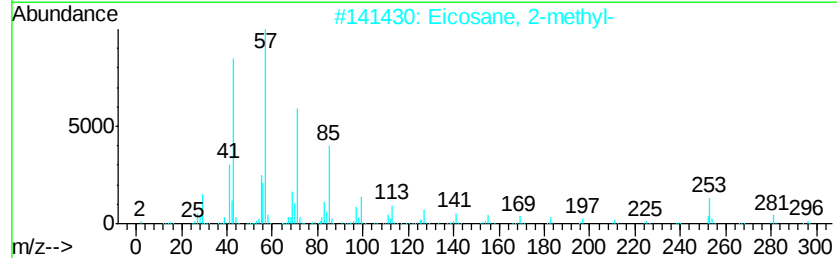
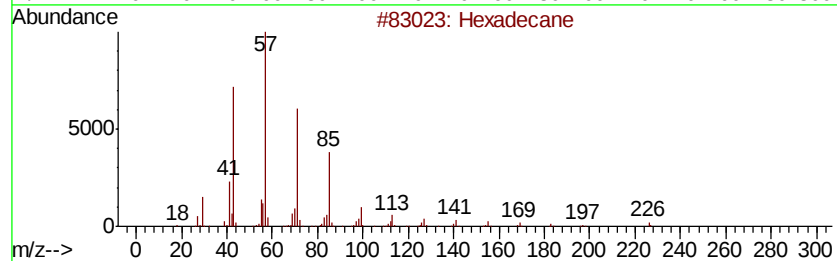
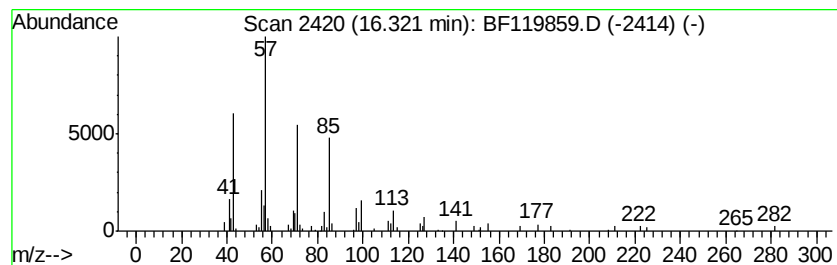
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Hexadecane Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.32 | 2.58 ng | 154398 | Perylene-d12     | 15.37 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#         | Qual |
|------|----|---------------------|-----|---------|--------------|------|
| 1    | 5  | Hexadecane          | 226 | C16H34  | 000544-76-3  | 64   |
| 2    |    | Eicosane, 2-methyl- | 296 | C21H44  | 001560-84-5  | 64   |
| 3    |    | Heptadecane         | 240 | C17H36  | 000629-78-7  | 64   |
| 4    |    | 2-methyloctacosane  | 408 | C29H60  | 1000376-72-8 | 64   |
| 5    |    | Heneicosane         | 296 | C21H44  | 000629-94-7  | 62   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\  
Data File : BF119859.D  
Acq On : 9 Apr 2020 7:38  
Operator : MA/SJ  
Sample : L2151-09  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MH-109

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

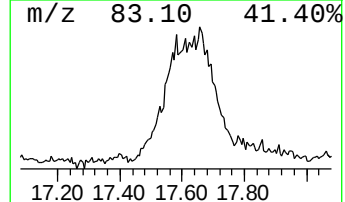
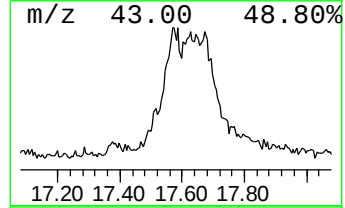
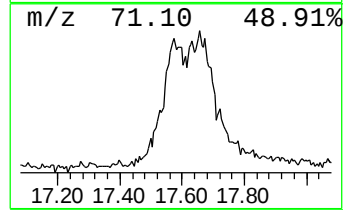
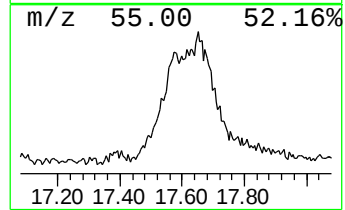
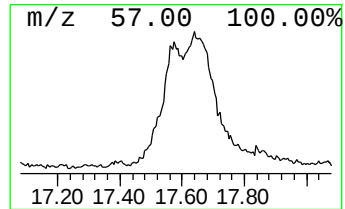
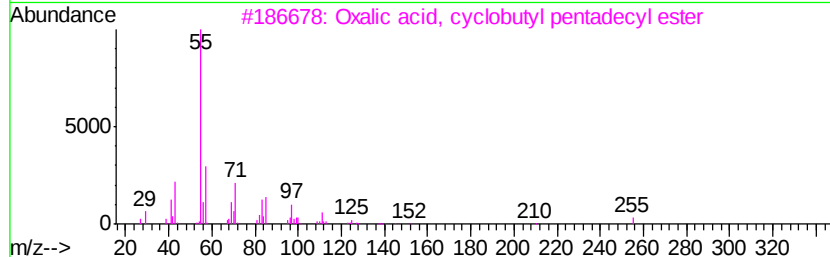
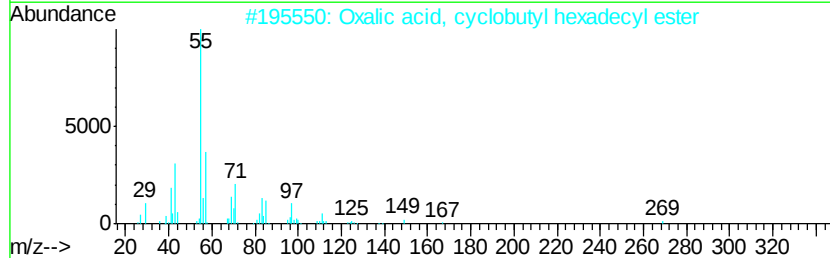
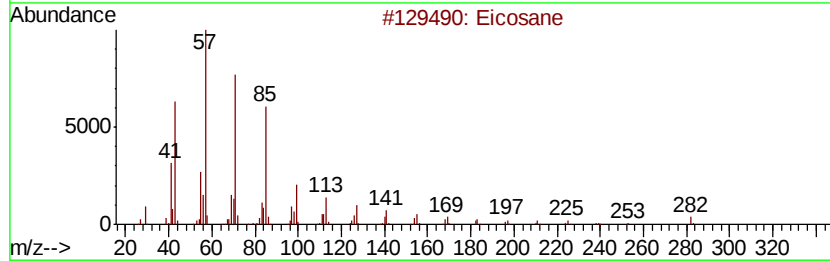
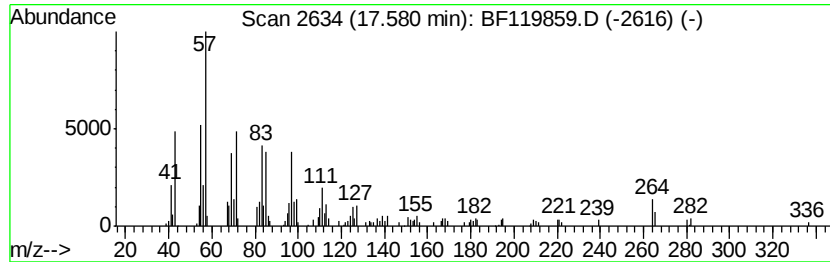
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Eicosane Concentration Rank 3

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.58 | 8.85 ng | 529765 | Perylene-d12     | 15.37 |

| Hit# | of | 5 | Tentative ID                             | MW  | MolForm    | CAS#         | Qual |
|------|----|---|------------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Eicosane                                 | 282 | C20H42     | 000112-95-8  | 95   |
| 2    |    |   | Oxalic acid, cyclobutyl hexadecyl ester  | 368 | C22H40O4   | 1000309-70-6 | 83   |
| 3    |    |   | Oxalic acid, cyclobutyl pentadecyl ester | 354 | C21H38O4   | 1000309-70-5 | 83   |
| 4    |    |   | Oxalic acid, isobutyl octadecyl ester    | 398 | C24H46O4   | 1000309-38-3 | 83   |
| 5    |    |   | Dotriacontyl heptafluorobutyrate         | 662 | C36H65F7O2 | 1000351-84-2 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF040820\  
Data File : BF119859.D  
Acq On : 9 Apr 2020 7:38  
Operator : MA/SJ  
Sample : L2151-09  
Misc :  
ALS Vial : 27 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MH-109

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF040820.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 |
| 2-Pentanone, 4-hy... | 4.96  | 2.3     | ng    | 119962   | 1                     | 6.76  | 1068640 | 20.0 |
| n-Hexadecanoic acid  | 11.83 | 7.8     | ng    | 728862   | 4                     | 11.29 | 1871270 | 20.0 |
| Octadecanoic acid    | 12.61 | 3.2     | ng    | 300338   | 4                     | 11.29 | 1871270 | 20.0 |
| Tricosyl pentaflu... | 13.79 | 9.0     | ng    | 673534   | 5                     | 13.93 | 1497510 | 20.0 |
| Octacosane           | 14.41 | 2.4     | ng    | 180065   | 5                     | 13.93 | 1497510 | 20.0 |
| Heneicosane          | 14.71 | 3.6     | ng    | 213834   | 6                     | 15.37 | 1197280 | 20.0 |
| Tetracosane          | 15.04 | 3.2     | ng    | 194157   | 6                     | 15.37 | 1197280 | 20.0 |
| 2-methyloctacosane   | 15.83 | 2.7     | ng    | 161162   | 6                     | 15.37 | 1197280 | 20.0 |
| Hexadecane           | 16.32 | 2.6     | ng    | 154398   | 6                     | 15.37 | 1197280 | 20.0 |
| Eicosane             | 17.58 | 8.8     | ng    | 529765   | 6                     | 15.37 | 1197280 | 20.0 |