

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\  
 Data File : BF115510.D  
 Acq On : 11 Jul 2019 22:28  
 Operator : HP/JU  
 Sample : K3751-03  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 DUCT-BANK-TEST-PIT

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-BF070519.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         | 2.187       | 11            | 17          | 33           | rVB      | 1277819        | 1736188       | 50.01%          | 5.253%        |
| 2         | 5.510       | 577           | 582         | 585          | rBV      | 2664404        | 2565633       | 73.90%          | 7.762%        |
| 3         | 6.516       | 748           | 753         | 756          | rBV      | 2778923        | 2742345       | 78.99%          | 8.297%        |
| 4         | 6.663       | 773           | 778         | 781          | rBV      | 3082233        | 2818021       | 81.17%          | 8.526%        |
| 5         | 6.892       | 813           | 817         | 820          | rBV      | 1137788        | 880631        | 25.37%          | 2.664%        |
| 6         | 7.051       | 840           | 844         | 847          | rBV      | 2689333        | 2185458       | 62.95%          | 6.612%        |
| 7         | 7.451       | 907           | 912         | 915          | rBV      | 2004734        | 1737021       | 50.03%          | 5.255%        |
| 8         | 8.175       | 1031          | 1035        | 1039         | rBV      | 1477768        | 1138222       | 32.79%          | 3.444%        |
| 9         | 9.251       | 1214          | 1218        | 1221         | rBV      | 3999898        | 3289068       | 94.74%          | 9.951%        |
| 10        | 9.639       | 1281          | 1284        | 1294         | rBV      | 870979         | 703189        | 20.25%          | 2.127%        |
| 11        | 9.933       | 1330          | 1334        | 1337         | rBV      | 1426878        | 1348622       | 38.85%          | 4.080%        |
| 12        | 10.716      | 1463          | 1467        | 1480         | rBV      | 2421233        | 2198168       | 63.32%          | 6.650%        |
| 13        | 11.416      | 1582          | 1586        | 1589         | rBV      | 1731500        | 1433177       | 41.28%          | 4.336%        |
| 14        | 11.945      | 1672          | 1676        | 1692         | rBV      | 511594         | 580732        | 16.73%          | 1.757%        |
| 15        | 12.727      | 1805          | 1809        | 1823         | rBV      | 275571         | 295832        | 8.52%           | 0.895%        |
| 16        | 13.004      | 1851          | 1856        | 1859         | rBV      | 3667483        | 3471711       | 100.00%         | 10.504%       |
| 17        | 13.898      | 2004          | 2008        | 2023         | rBV2     | 205649         | 347314        | 10.00%          | 1.051%        |
| 18        | 14.051      | 2030          | 2034        | 2038         | rBV      | 1162640        | 1105488       | 31.84%          | 3.345%        |
| 19        | 14.221      | 2060          | 2063        | 2071         | rBV      | 63019          | 65825         | 1.90%           | 0.199%        |
| 20        | 14.527      | 2113          | 2115        | 2121         | rVB      | 104065         | 88819         | 2.56%           | 0.269%        |
| 21        | 14.839      | 2164          | 2168        | 2171         | rBV      | 150895         | 149644        | 4.31%           | 0.453%        |
| 22        | 14.915      | 2178          | 2181        | 2188         | rVB      | 34569          | 40948         | 1.18%           | 0.124%        |
| 23        | 15.186      | 2222          | 2227        | 2237         | rBV      | 185589         | 242462        | 6.98%           | 0.734%        |
| 24        | 15.533      | 2280          | 2286        | 2289         | rBV      | 877562         | 1148178       | 33.07%          | 3.474%        |
| 25        | 15.574      | 2289          | 2293        | 2299         | rVB      | 205472         | 287496        | 8.28%           | 0.870%        |
| 26        | 16.015      | 2363          | 2368        | 2373         | rBV      | 152735         | 244041        | 7.03%           | 0.738%        |
| 27        | 16.533      | 2451          | 2456        | 2461         | rBV      | 81986          | 139200        | 4.01%           | 0.421%        |
| 28        | 17.139      | 2555          | 2559        | 2567         | rVB3     | 31691          | 69267         | 2.00%           | 0.210%        |

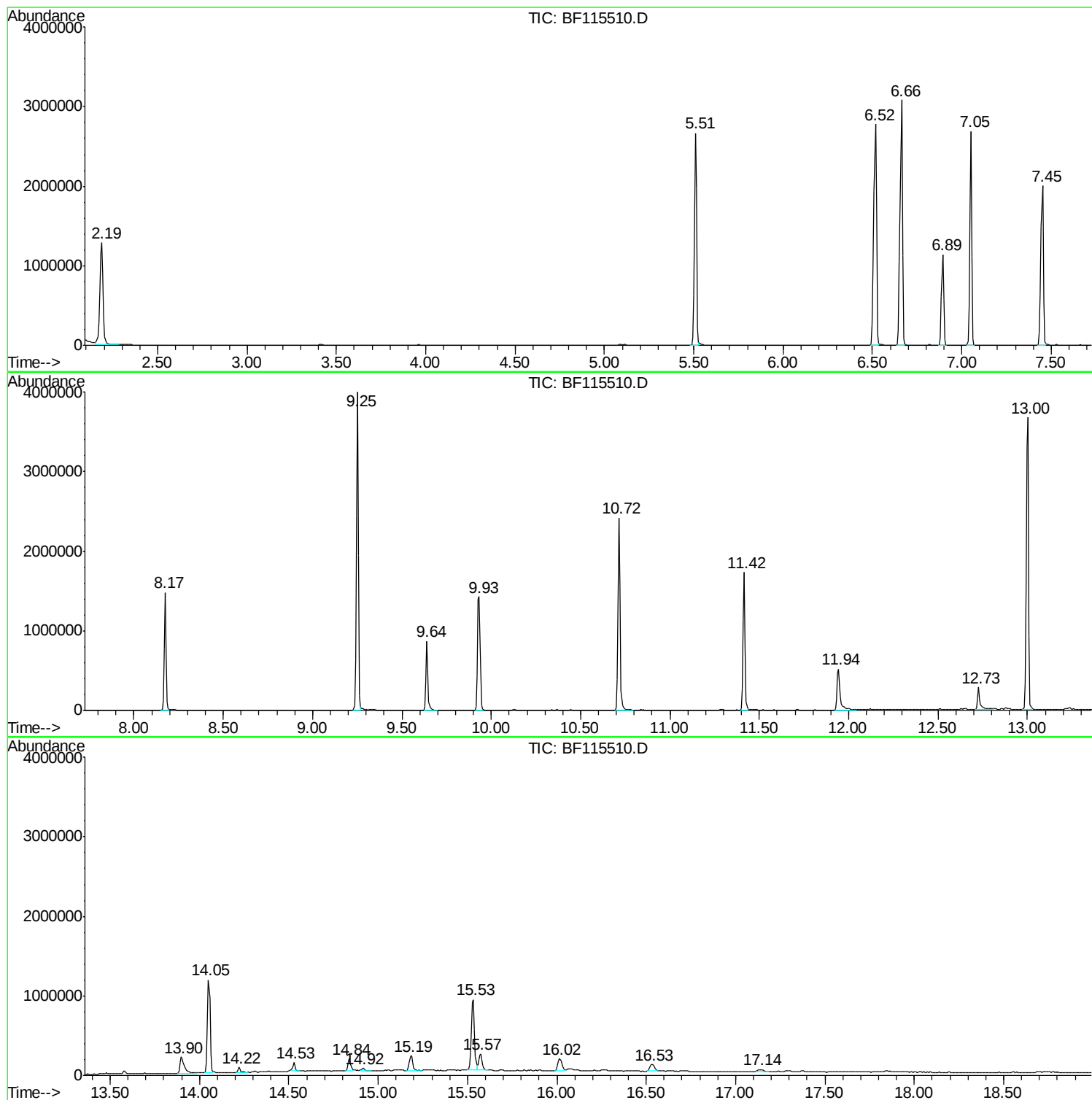
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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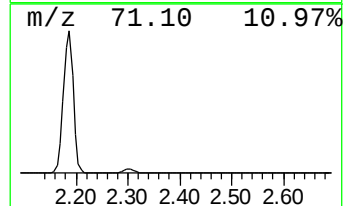
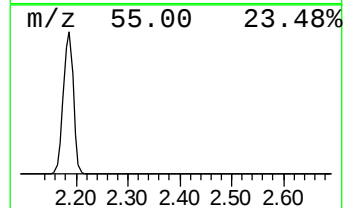
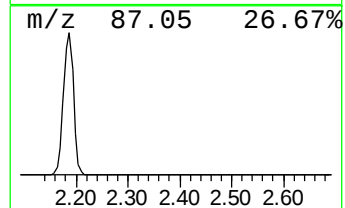
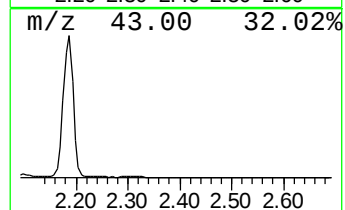
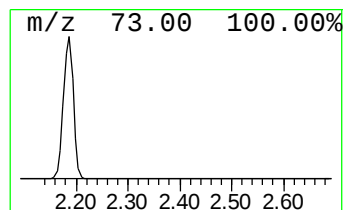
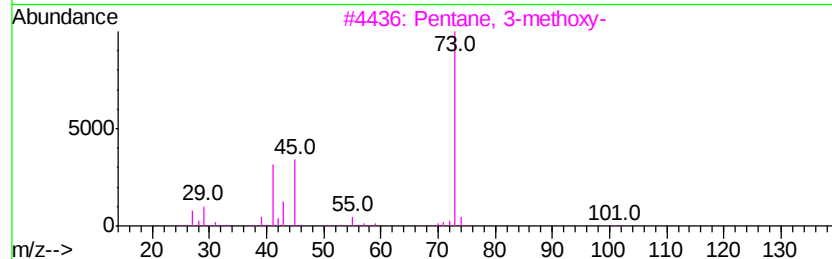
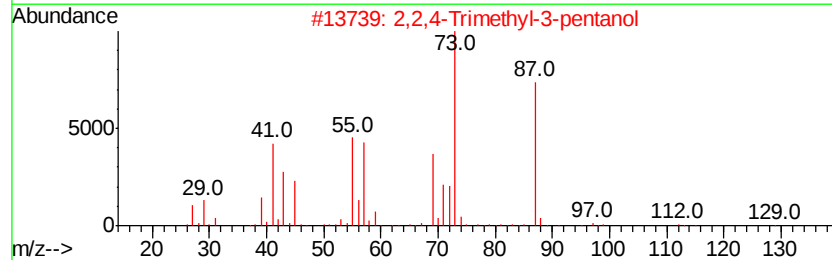
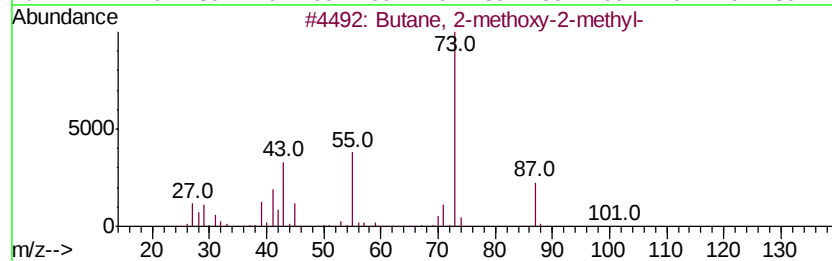
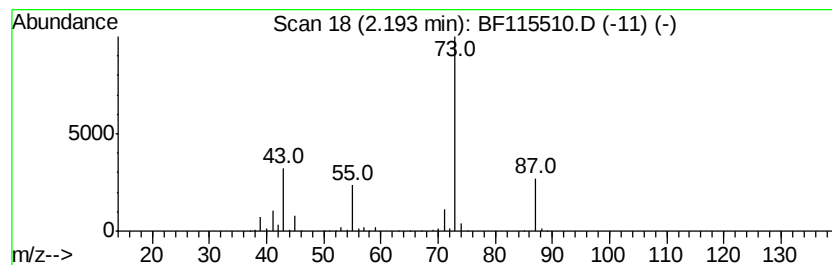
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T.    | EstConc  | Area                        | Relative to ISTD       | R.T.    |             |      |
|---------|----------|-----------------------------|------------------------|---------|-------------|------|
| 2.19    | 39.43 ng | 1736190                     | 1,4-Dichlorobenzene-d4 | 6.89    |             |      |
| Hit# of | 5        | Tentative ID                | MW                     | MolForm | CAS#        | Qual |
| 1       |          | 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       |          | Pentane, 3-methoxy-         | 102                    | C6H14O  | 036839-67-5 | 23   |
| 4       |          | Silane, tetramethyl-        | 88                     | C4H12Si | 000075-76-3 | 17   |
| 5       |          | N-Ethylformamide            | 73                     | C3H7NO  | 000627-45-2 | 9    |



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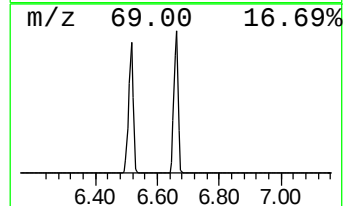
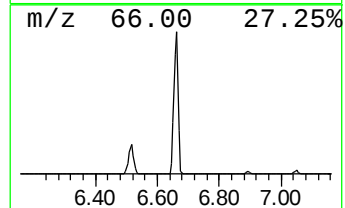
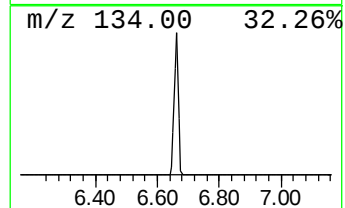
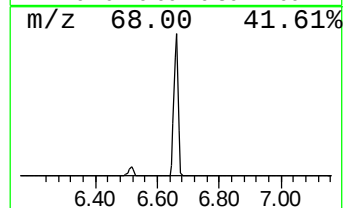
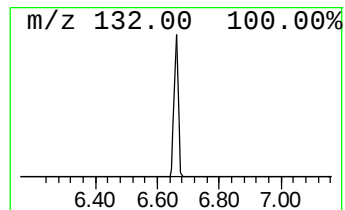
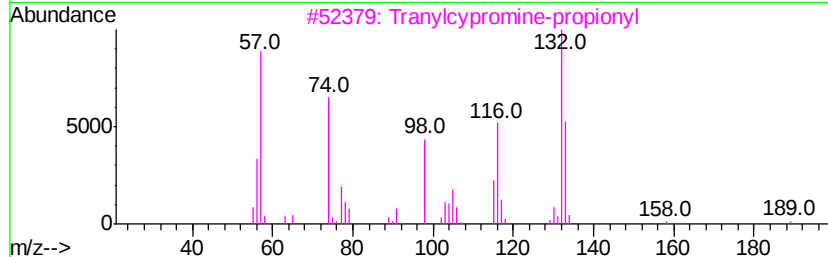
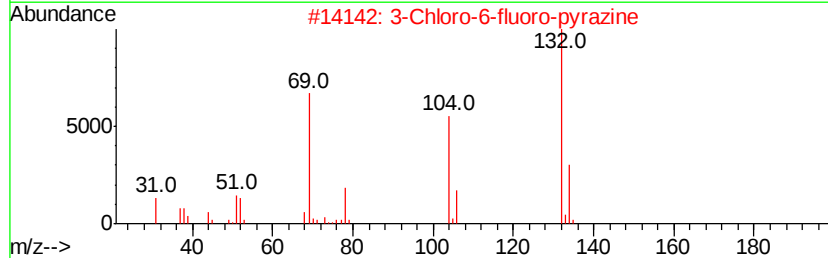
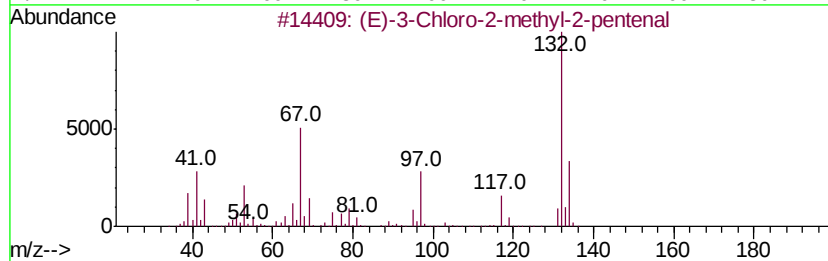
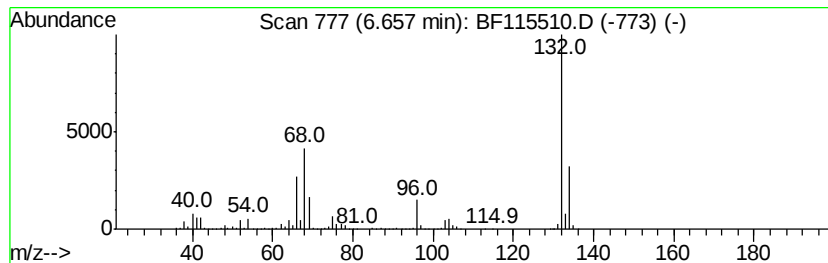
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown6.66 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.66 | 64.00 ng | 2818020 | 1,4-Dichlorobenzene-d4 | 6.89 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 2    |    | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 16   |
| 3    |    | Tranvlcypropine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 4-Methylpyrrolo[1,2-a]pyrazine      | 132 | C8H8N2    | 064608-60-2  | 9    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

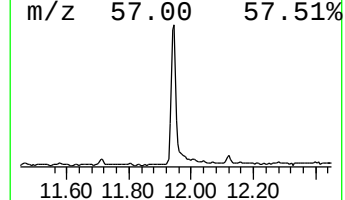
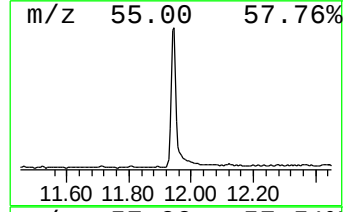
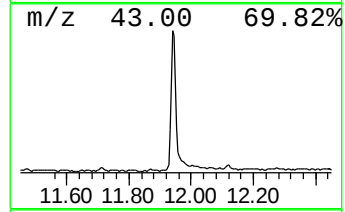
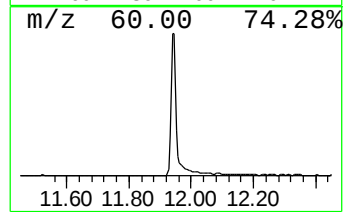
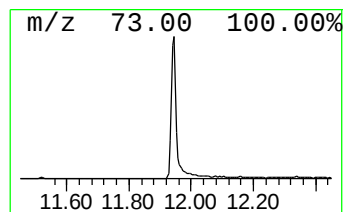
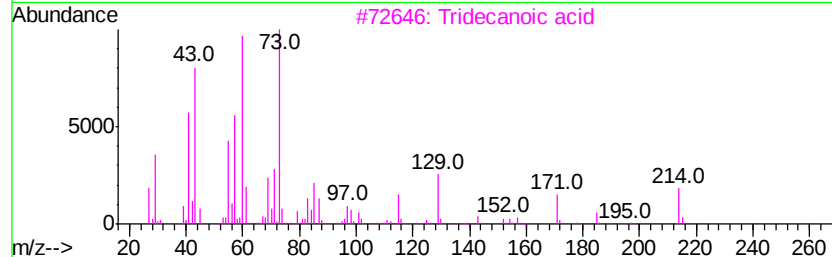
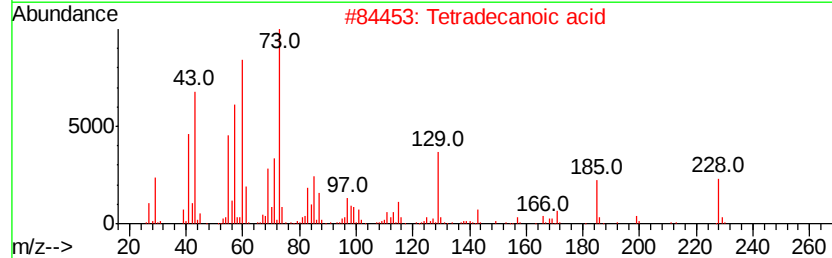
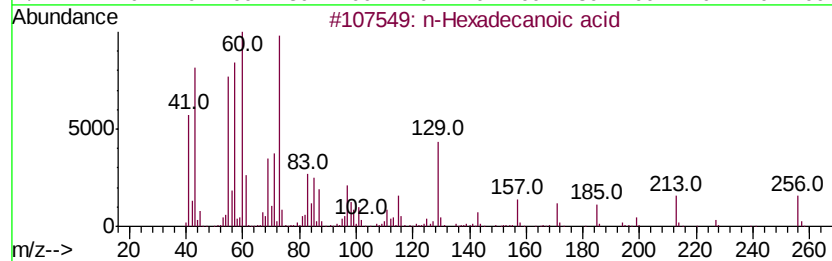
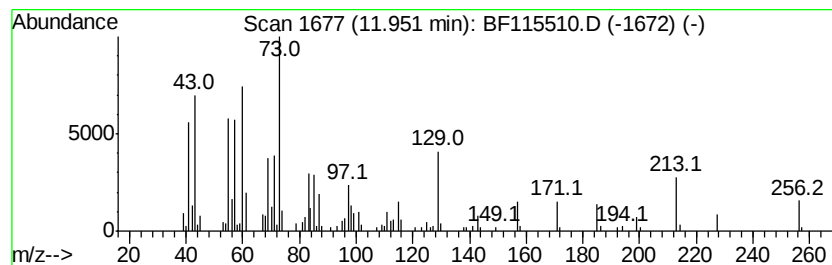
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.95 | 8.10 ng | 580732 | Phenanthrene-d10 | 11.42 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 2    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 94   |
| 3    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 91   |
| 4    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 83   |
| 5    |    | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 76   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M  
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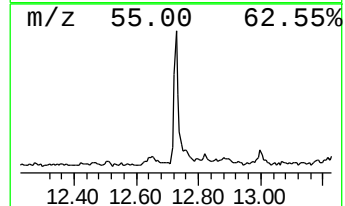
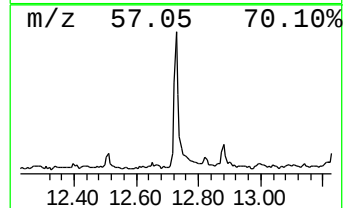
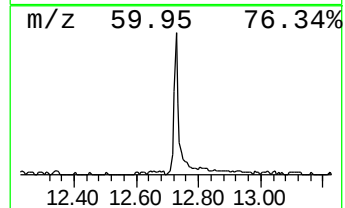
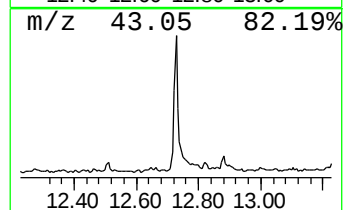
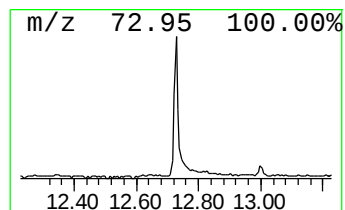
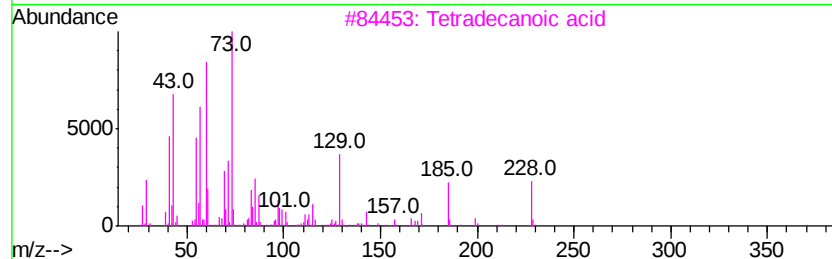
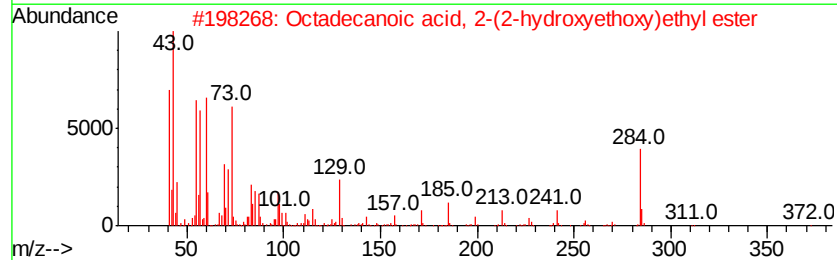
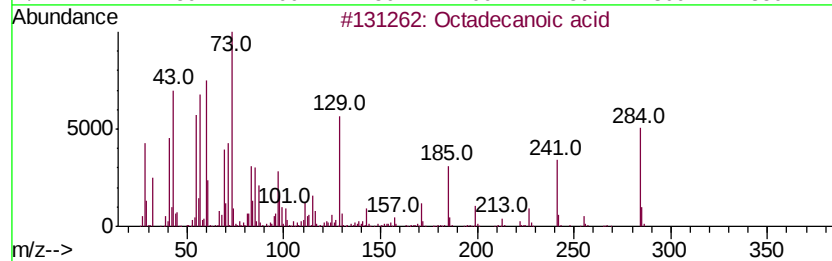
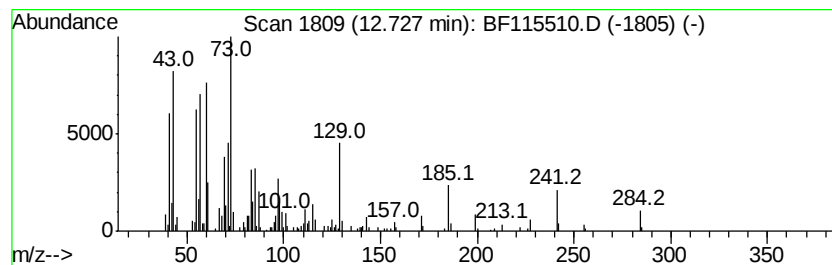
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Octadecanoic acid Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.73 | 4.13 ng | 295832 | Phenanthrene-d10 | 11.42 |

| Hit# of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|---------|---|--------------------------------------|-----|----------|-------------|------|
| 1       |   | Octadecanoic acid                    | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2       |   | Octadecanoic acid, 2-(2-hydroxyve... | 372 | C22H44O4 | 000106-11-6 | 91   |
| 3       |   | Tetradecanoic acid                   | 228 | C14H28O2 | 000544-63-8 | 83   |
| 4       |   | Pentadecanoic acid                   | 242 | C15H30O2 | 001002-84-2 | 81   |
| 5       |   | n-Decanoic acid                      | 172 | C10H20O2 | 000334-48-5 | 70   |



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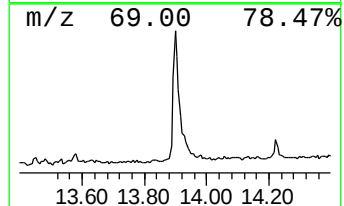
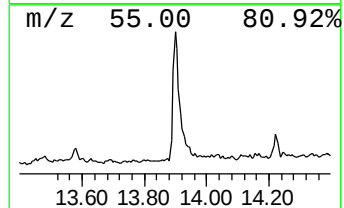
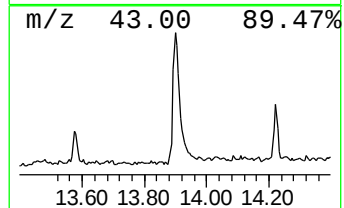
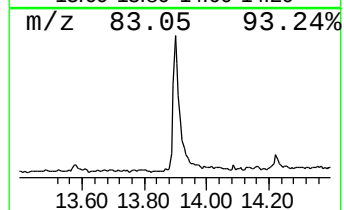
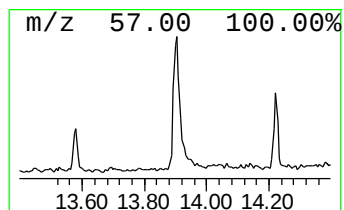
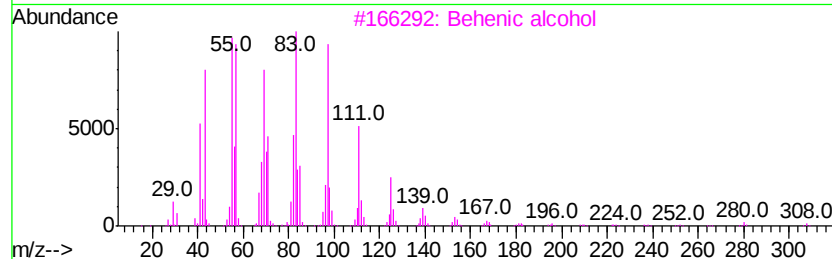
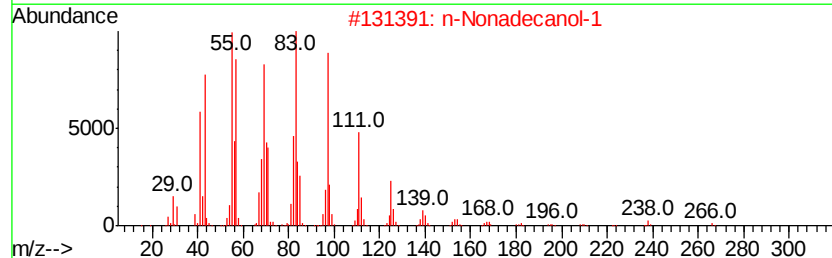
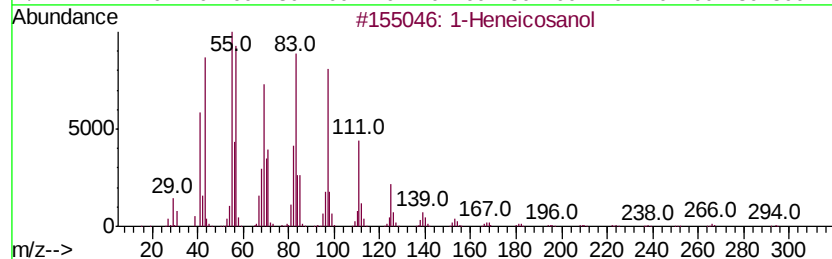
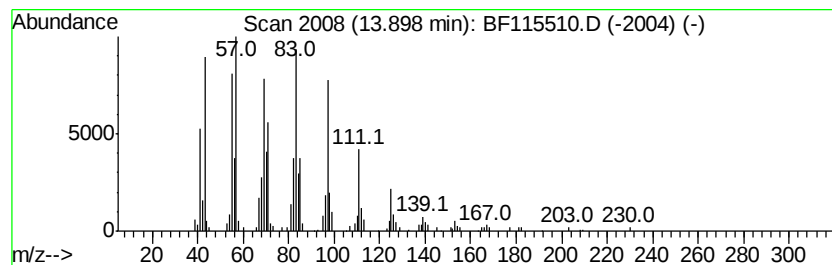
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 1-Heneicosanol Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.90 | 6.28 ng | 347314 | Chrysene-d12     | 14.05 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|---------|---|-------------------------------------|-----|------------|-------------|------|
| 1       |   | 1-Heneicosanol                      | 312 | C21H44O    | 015594-90-8 | 95   |
| 2       |   | n-Nonadecanol-1                     | 284 | C19H40O    | 001454-84-8 | 95   |
| 3       |   | Behenic alcohol                     | 326 | C22H46O    | 000661-19-8 | 95   |
| 4       |   | 1-Heptacosanol                      | 396 | C27H56O    | 002004-39-9 | 94   |
| 5       |   | Heptafluorobutyric acid, pentade... | 424 | C19H31F7O2 | 959261-23-5 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\  
 Data File : BF115510.D  
 Acq On : 11 Jul 2019 22:28  
 Operator : HP/JU  
 Sample : K3751-03  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 DUCT-BANK-TEST-PIT

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

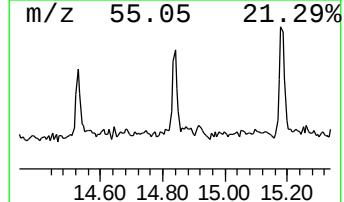
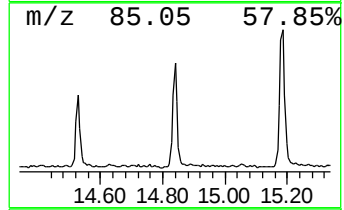
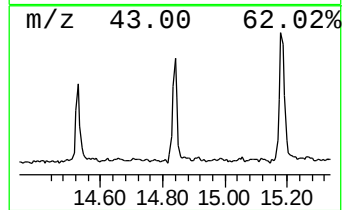
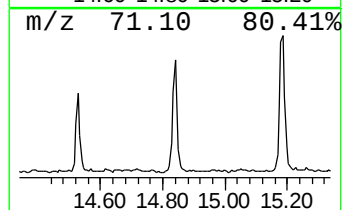
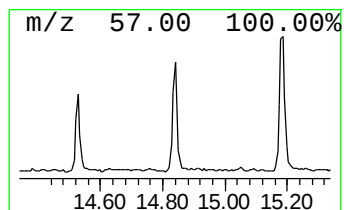
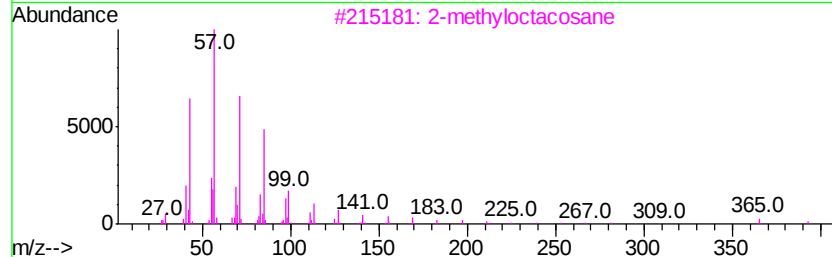
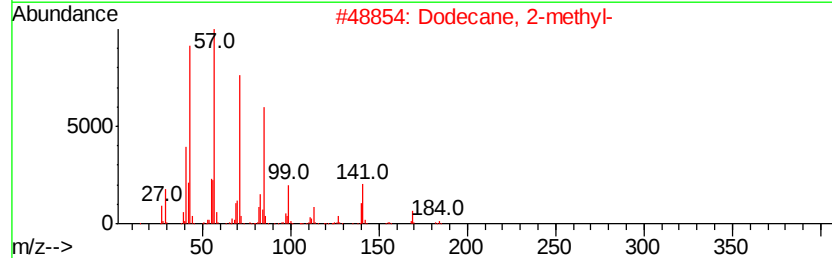
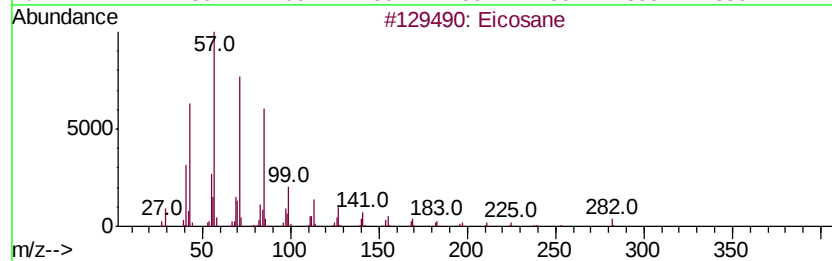
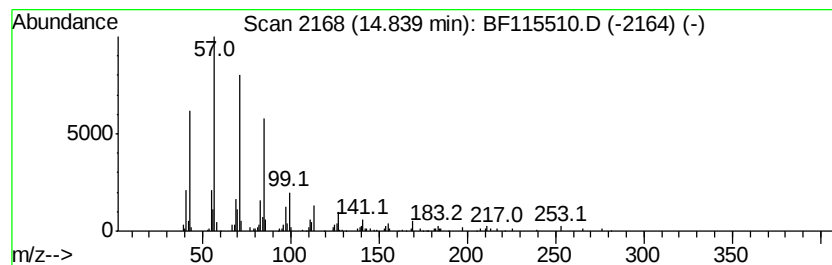
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 7 Eicosane Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.84 | 2.61 ng | 149644 | Perylene-d12     | 15.53 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#         | Qual |
|------|----|---------------------|-----|---------|--------------|------|
| 1    | 5  | Eicosane            | 282 | C20H42  | 000112-95-8  | 95   |
| 2    |    | Dodecane, 2-methyl- | 184 | C13H28  | 001560-97-0  | 90   |
| 3    |    | 2-methyloctacosane  | 408 | C29H60  | 1000376-72-8 | 90   |
| 4    |    | Hexacosane          | 366 | C26H54  | 000630-01-3  | 87   |
| 5    |    | Heneicosane         | 296 | C21H44  | 000629-94-7  | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\  
Data File : BF115510.D  
Acq On : 11 Jul 2019 22:28  
Operator : HP/JU  
Sample : K3751-03  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

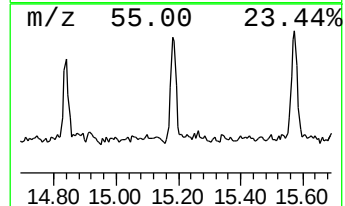
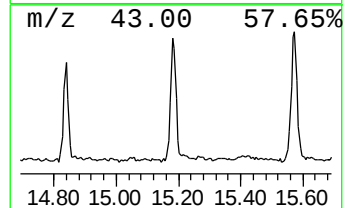
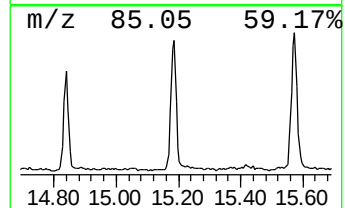
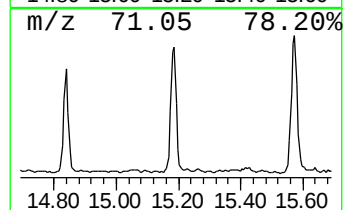
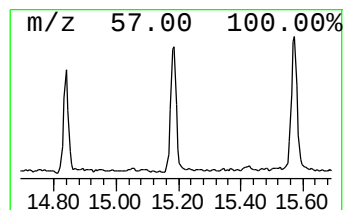
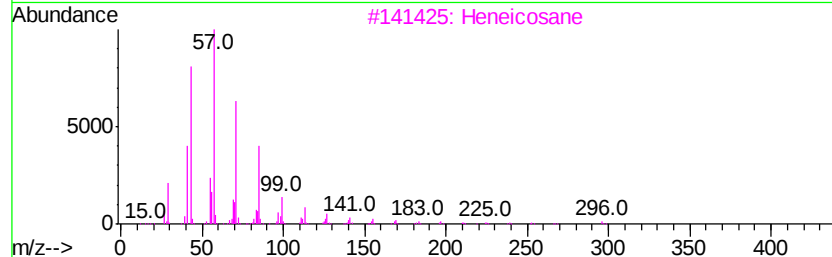
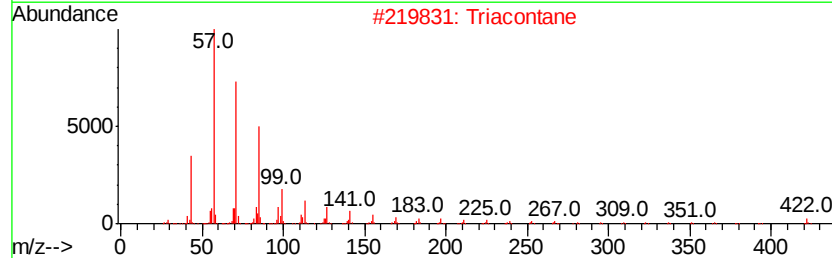
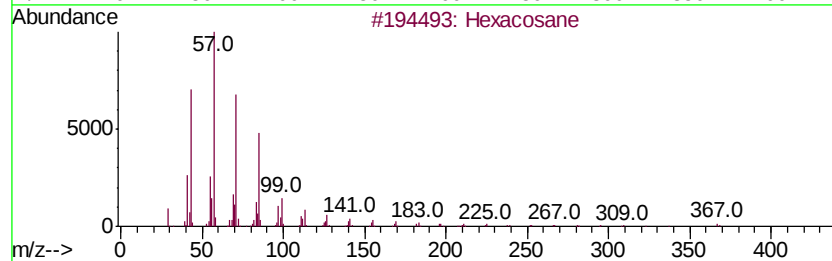
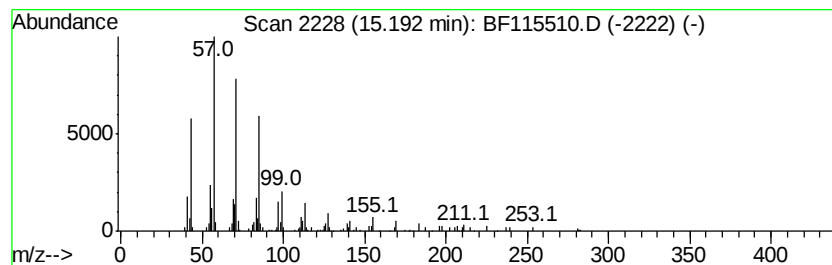
Instrument :  
BNA\_F  
ClientSampleId :  
DUCT-BANK-TEST-PIT

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.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 Hexacosane Concentration Rank 7

| R.T.    | EstConc               | Area         | Relative to ISTD | R.T.      |
|---------|-----------------------|--------------|------------------|-----------|
| 15.19   | 4.22 ng               | 242462       | Perylene-d12     | 15.53     |
| Hit# of | 5                     | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Hexacosane            | 366 C26H54   | 000630-01-3      | 91        |
| 2       | triacontane           | 422 C30H62   | 000638-68-6      | 91        |
| 3       | Heneicosane           | 296 C21H44   | 000629-94-7      | 91        |
| 4       | 2-methyloctacosane    | 408 C29H60   | 1000376-72-8     | 91        |
| 5       | Heptadecane, 9-octyl- | 352 C25H52   | 007225-64-1      | 91        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\  
Data File : BF115510.D  
Acq On : 11 Jul 2019 22:28  
Operator : HP/JU  
Sample : K3751-03  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DUCT-BANK-TEST-PIT

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

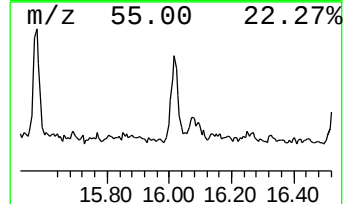
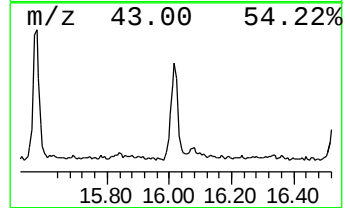
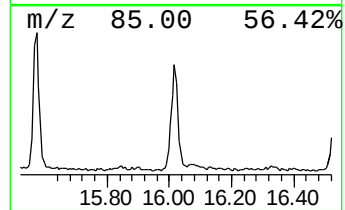
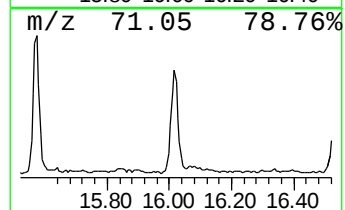
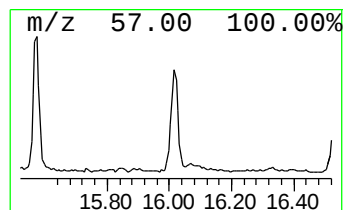
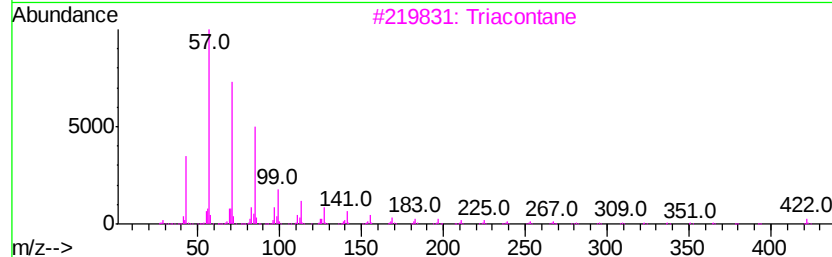
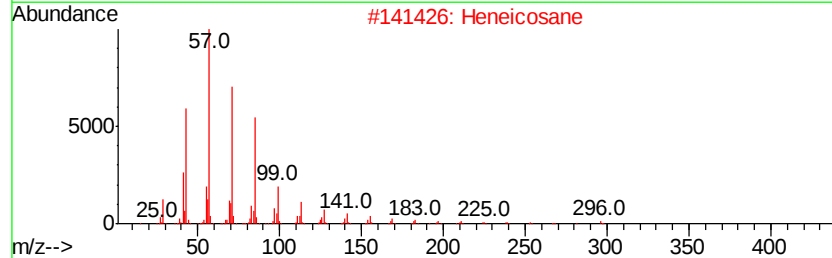
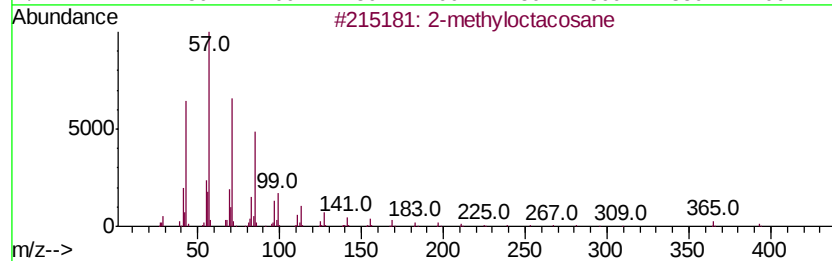
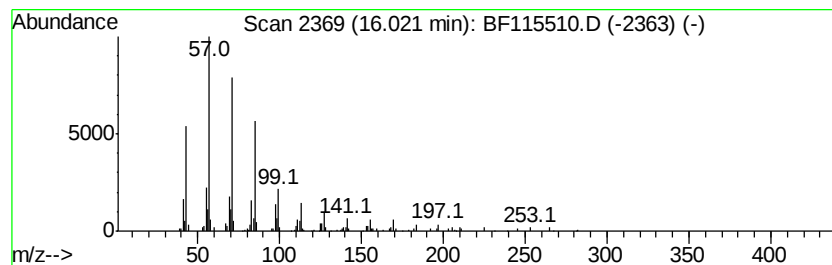
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 2-methyloctacosane Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.02 | 4.25 ng | 244041 | Perylene-d12     | 15.53 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------|-----|---------|--------------|------|
| 1    | 5  | 2-methyloctacosane | 408 | C29H60  | 1000376-72-8 | 91   |
| 2    |    | Heneicosane        | 296 | C21H44  | 000629-94-7  | 91   |
| 3    |    | Triacotane         | 422 | C30H62  | 000638-68-6  | 91   |
| 4    |    | Heptacosane        | 380 | C27H56  | 000593-49-7  | 91   |
| 5    |    | Hexacosane         | 366 | C26H54  | 000630-01-3  | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\  
Data File : BF115510.D  
Acq On : 11 Jul 2019 22:28  
Operator : HP/JU  
Sample : K3751-03  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DUCT-BANK-TEST-PIT

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

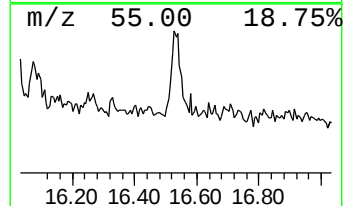
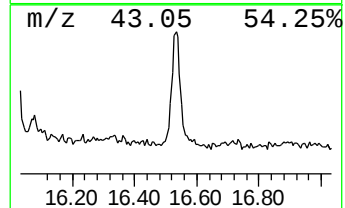
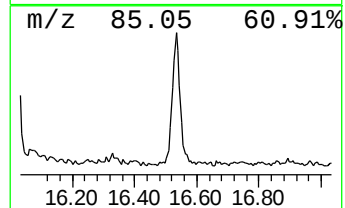
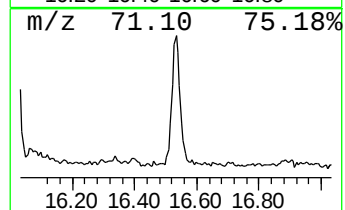
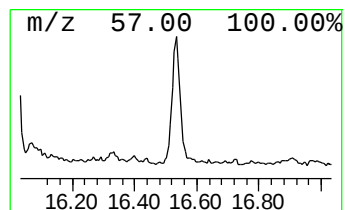
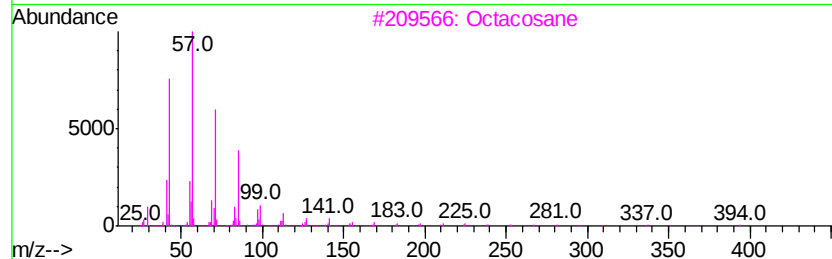
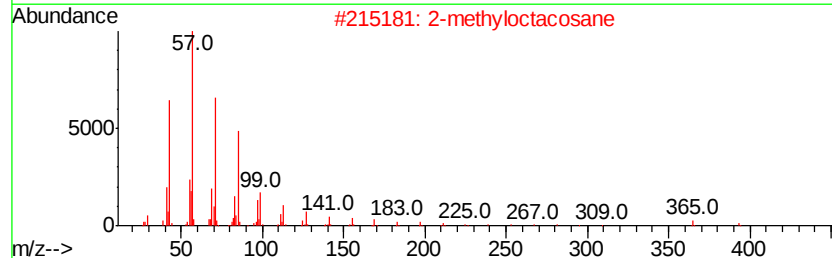
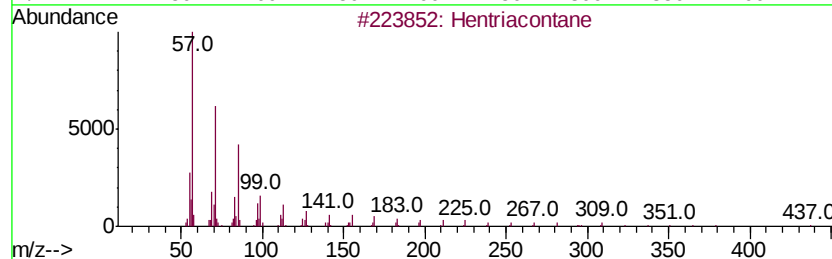
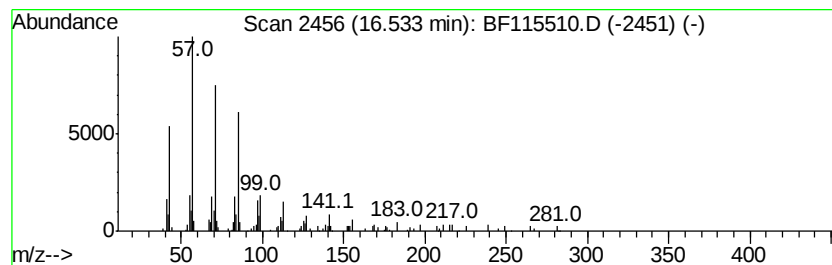
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Hentriacontane Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.53 | 2.42 ng | 139200 | Perylene-d12     | 15.53 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm | CAS#         | Qual |
|------|----|---|--------------------|-----|---------|--------------|------|
| 1    |    |   | Hentriacontane     | 437 | C31H64  | 000630-04-6  | 91   |
| 2    |    |   | 2-methyloctacosane | 408 | C29H60  | 1000376-72-8 | 91   |
| 3    |    |   | Octacosane         | 394 | C28H58  | 000630-02-4  | 91   |
| 4    |    |   | Triacosane         | 422 | C30H62  | 000638-68-6  | 91   |
| 5    |    |   | Heptacosane        | 380 | C27H56  | 000593-49-7  | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF071119\  
Data File : BF115510.D  
Acq On : 11 Jul 2019 22:28  
Operator : HP/JU  
Sample : K3751-03  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
DUCT-BANK-TEST-PIT

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF070519.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... | 2.19  | 39.4    | ng    | 1736190  | 1                     | 6.89  | 880631  | 20.0 |
| unknown6.66          | 6.66  | 64.0    | ng    | 2818020  | 1                     | 6.89  | 880631  | 20.0 |
| n-Hexadecanoic acid  | 11.95 | 8.1     | ng    | 580732   | 4                     | 11.42 | 1433180 | 20.0 |
| Octadecanoic acid    | 12.73 | 4.1     | ng    | 295832   | 4                     | 11.42 | 1433180 | 20.0 |
| 1-Heneicosanol       | 13.90 | 6.3     | ng    | 347314   | 5                     | 14.05 | 1105490 | 20.0 |
| Eicosane             | 14.84 | 2.6     | ng    | 149644   | 6                     | 15.53 | 1148180 | 20.0 |
| Hexacosane           | 15.19 | 4.2     | ng    | 242462   | 6                     | 15.53 | 1148180 | 20.0 |
| 2-methyloctacosane   | 16.02 | 4.3     | ng    | 244041   | 6                     | 15.53 | 1148180 | 20.0 |
| Hentriacontane       | 16.53 | 2.4     | ng    | 139200   | 6                     | 15.53 | 1148180 | 20.0 |