

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF120319\  
 Data File : BF118061.D  
 Acq On : 3 Dec 2019 14:37  
 Operator : JU  
 Sample : PB125174BL  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB125174BL

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-BF111319.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         | 2.152       | 4             | 11          | 16           | rVB      | 3746526        | 5005516       | 91.30%          | 12.158%       |
| 2         | 5.087       | 506           | 510         | 518          | rVB      | 293987         | 343702        | 6.27%           | 0.835%        |
| 3         | 5.492       | 574           | 579         | 582          | rBV      | 2566885        | 2242293       | 40.90%          | 5.446%        |
| 4         | 6.498       | 746           | 750         | 753          | rBV      | 1637721        | 1443739       | 26.33%          | 3.507%        |
| 5         | 6.645       | 771           | 775         | 778          | rBV      | 4054276        | 4005257       | 73.05%          | 9.729%        |
| 6         | 6.875       | 811           | 814         | 818          | rBV      | 1029313        | 827700        | 15.10%          | 2.010%        |
| 7         | 7.034       | 837           | 841         | 844          | rBV      | 3765205        | 3196822       | 58.31%          | 7.765%        |
| 8         | 7.439       | 905           | 910         | 913          | rBV      | 2483766        | 2548881       | 46.49%          | 6.191%        |
| 9         | 8.163       | 1029          | 1033        | 1036         | rBV      | 1292508        | 1068578       | 19.49%          | 2.596%        |
| 10        | 9.245       | 1212          | 1217        | 1220         | rBV      | 5337378        | 4878997       | 88.99%          | 11.851%       |
| 11        | 9.839       | 1313          | 1318        | 1325         | rVB4     | 29824          | 60350         | 1.10%           | 0.147%        |
| 12        | 9.927       | 1329          | 1333        | 1336         | rBV      | 1468426        | 1256950       | 22.93%          | 3.053%        |
| 13        | 10.722      | 1463          | 1468        | 1471         | rBV      | 3399303        | 3191495       | 58.21%          | 7.752%        |
| 14        | 11.416      | 1583          | 1586        | 1590         | rBV      | 1463845        | 1328962       | 24.24%          | 3.228%        |
| 15        | 11.939      | 1668          | 1675        | 1680         | rBV      | 92099          | 126121        | 2.30%           | 0.306%        |
| 16        | 12.727      | 1805          | 1809        | 1821         | rBV2     | 95153          | 143864        | 2.62%           | 0.349%        |
| 17        | 13.016      | 1853          | 1858        | 1861         | rBV      | 5640183        | 5482705       | 100.00%         | 13.317%       |
| 18        | 14.068      | 2033          | 2037        | 2041         | rVB      | 1234671        | 1050486       | 19.16%          | 2.552%        |
| 19        | 14.227      | 2060          | 2064        | 2067         | rVB      | 91702          | 80550         | 1.47%           | 0.196%        |
| 20        | 14.533      | 2112          | 2116        | 2123         | rBV      | 201420         | 190166        | 3.47%           | 0.462%        |
| 21        | 14.845      | 2165          | 2169        | 2174         | rBV      | 305694         | 308875        | 5.63%           | 0.750%        |
| 22        | 15.192      | 2223          | 2228        | 2235         | rBV      | 308565         | 389167        | 7.10%           | 0.945%        |
| 23        | 15.562      | 2286          | 2291        | 2300         | rBV2     | 731385         | 1315353       | 23.99%          | 3.195%        |
| 24        | 16.039      | 2366          | 2372        | 2378         | rVB      | 214053         | 334101        | 6.09%           | 0.812%        |
| 25        | 16.556      | 2454          | 2460        | 2469         | rBV      | 123921         | 242902        | 4.43%           | 0.590%        |
| 26        | 17.174      | 2561          | 2565        | 2573         | rVB2     | 53176          | 106184        | 1.94%           | 0.258%        |

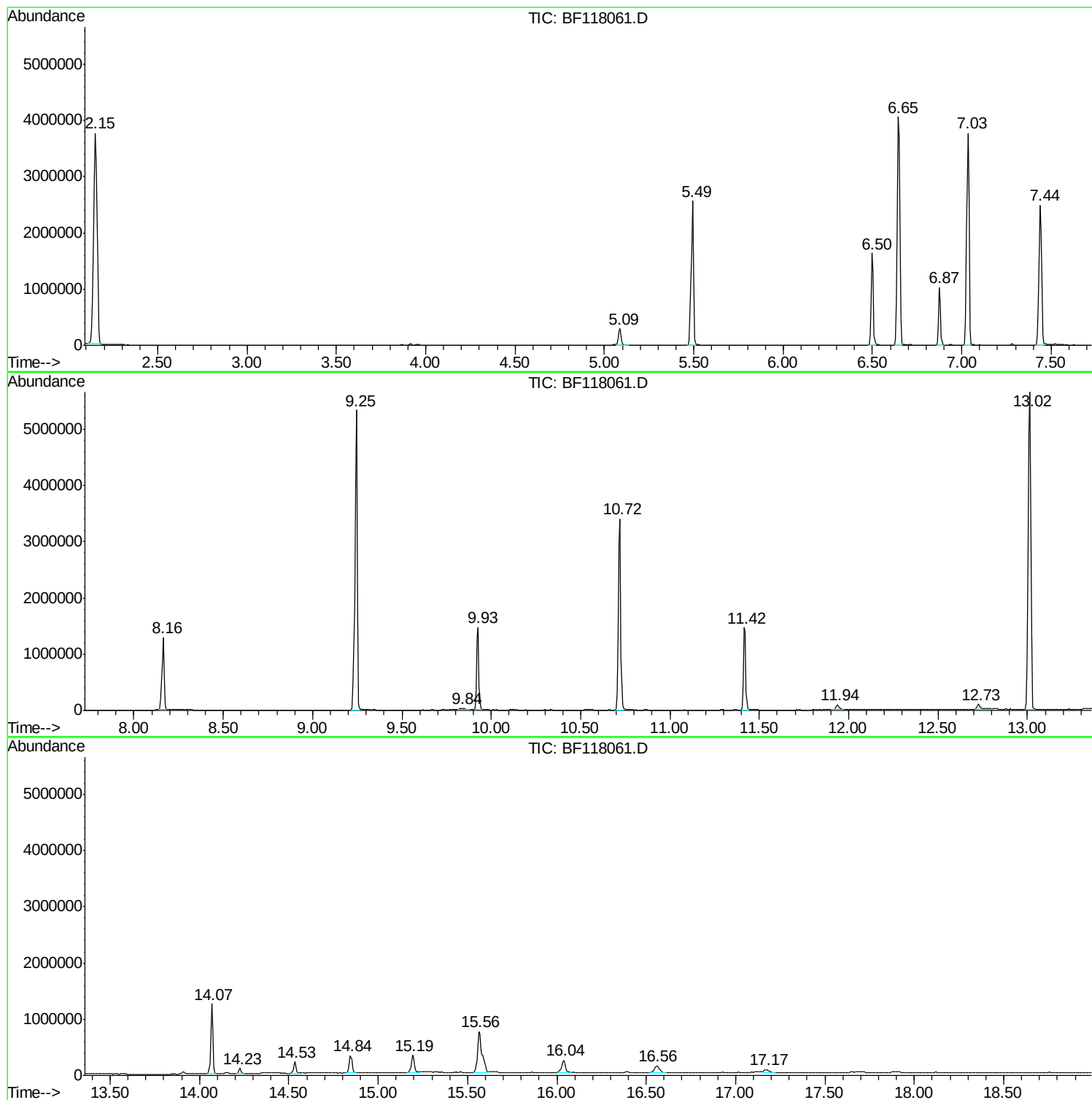
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ALS Vial : 4 Sample Multiplier: 1

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BNA\_F  
ClientSampleId :  
PB125174BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.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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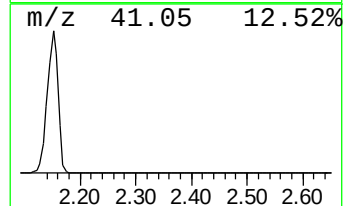
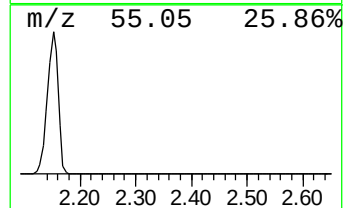
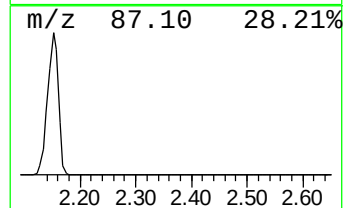
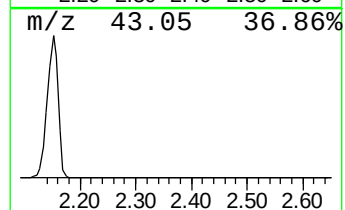
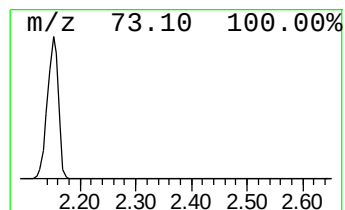
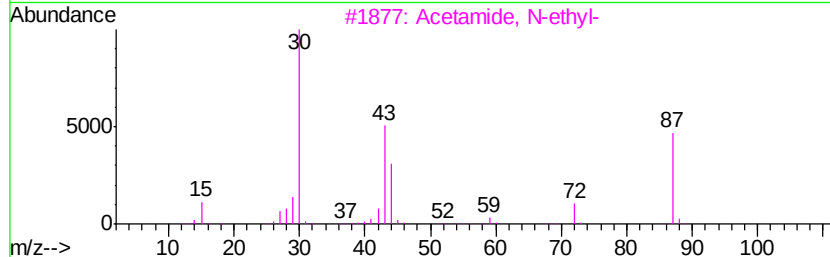
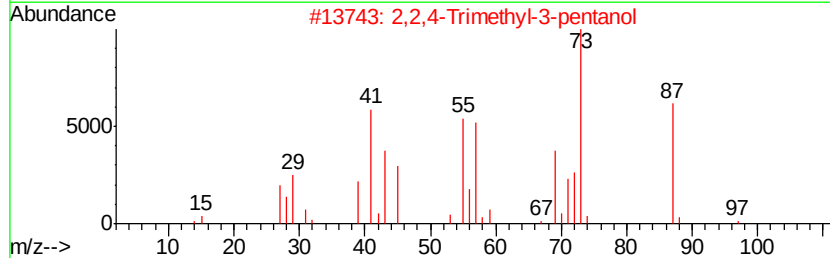
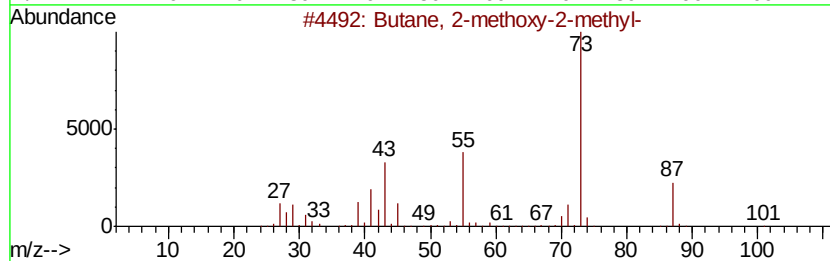
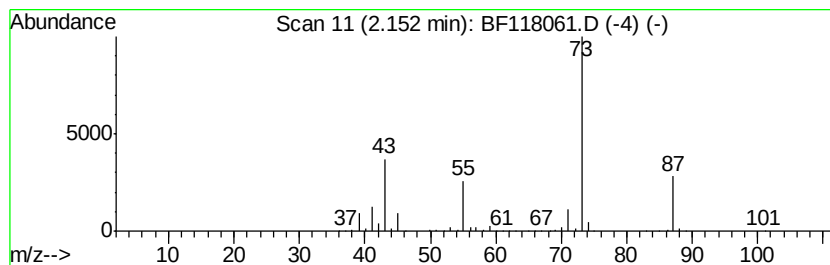
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 2.15 | 120.95 ng | 5005520 | 1,4-Dichlorobenzene-d4 | 6.87 |

| 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 | 40   |
| 3    |    | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 22   |
| 4    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 5    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 9    |



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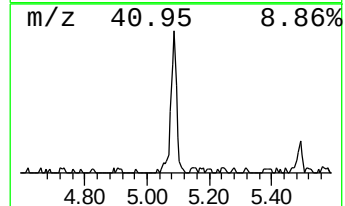
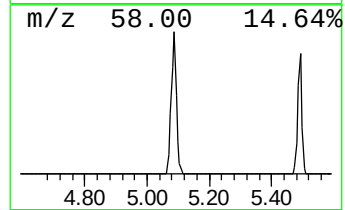
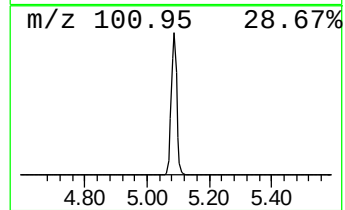
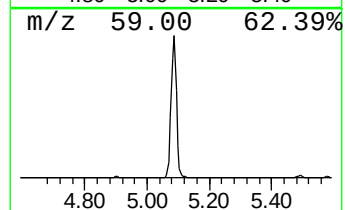
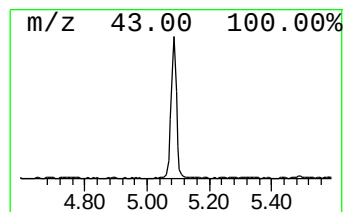
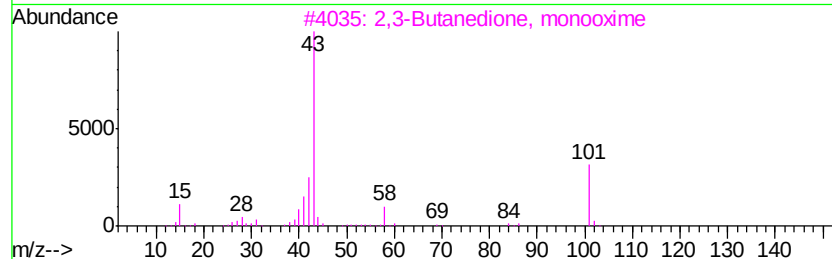
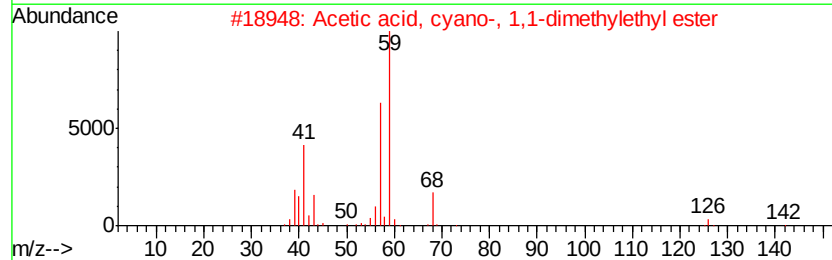
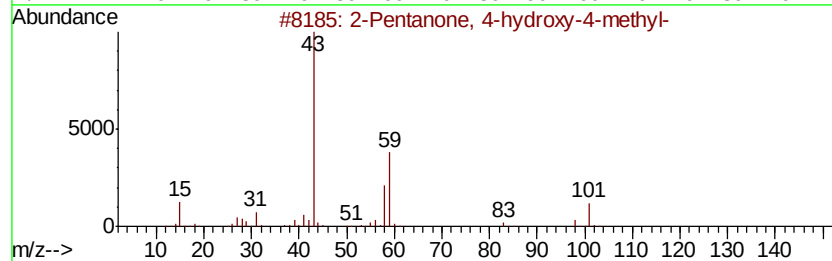
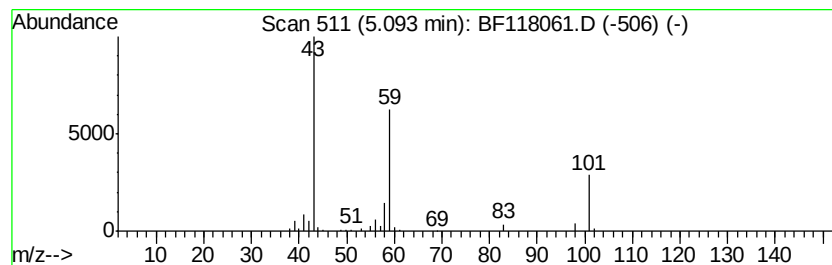
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.09 | 8.30 ng | 343702 | 1,4-Dichlorobenzene-d4 | 6.87 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-     | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    | Acetic acid, cyano-, 1,1-dimethyl... | 141 | C7H11NO2 | 001116-98-9 | 32   |
| 3    |    | 2,3-Butanedione, monooxime           | 101 | C4H7NO2  | 000057-71-6 | 17   |
| 4    |    | Morpholine, 4-methyl-                | 101 | C5H11NO  | 000109-02-4 | 9    |
| 5    |    | Butane, 1-ethoxy-                    | 102 | C6H14O   | 000628-81-9 | 9    |



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ALS Vial : 4 Sample Multiplier: 1

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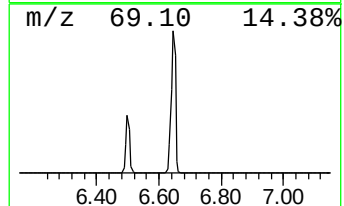
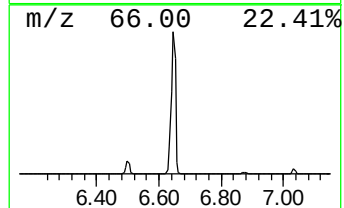
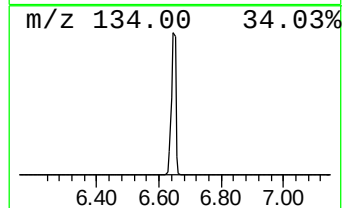
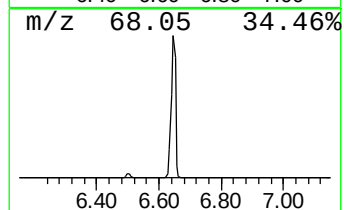
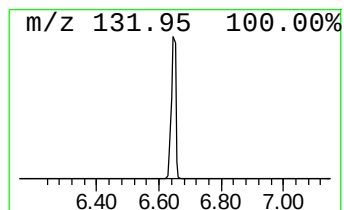
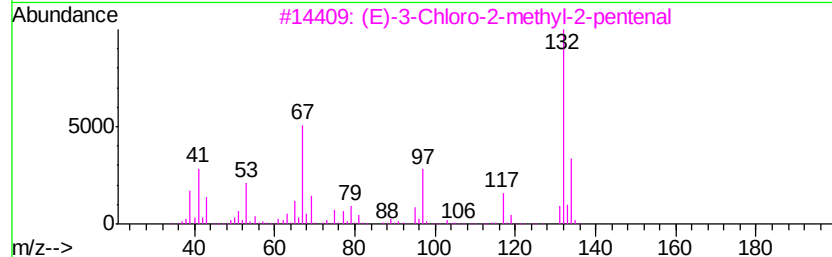
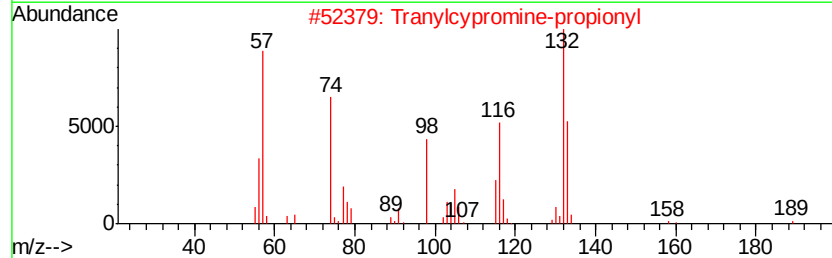
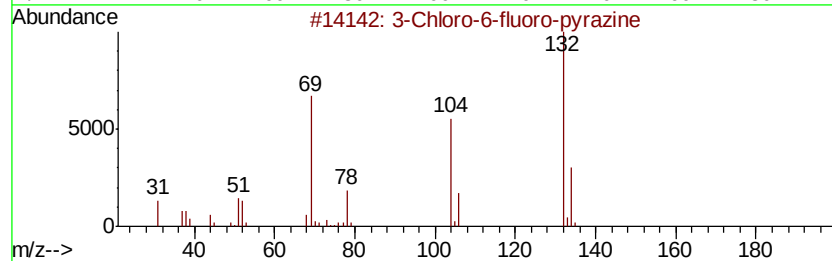
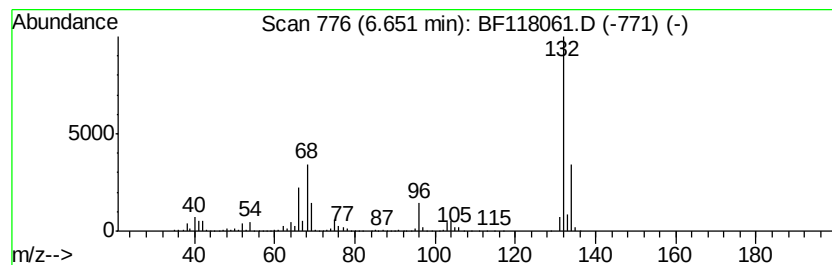
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 unknown6.65 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.65 | 96.78 ng | 4005260 | 1,4-Dichlorobenzene-d4 | 6.87 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    |   | Tranylcypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 3    |    |   | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 4    |    |   | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    |   | Benzene, 1-ethenyl-3,5-dimethyl-    | 132 | C10H12    | 005379-20-4  | 9    |



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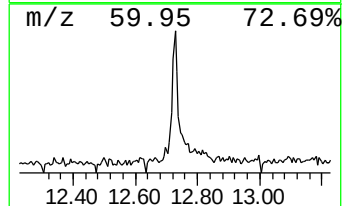
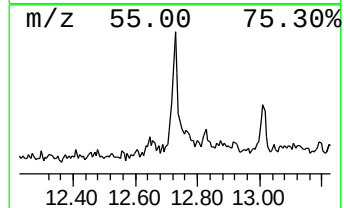
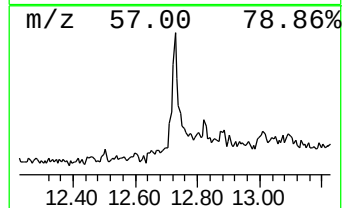
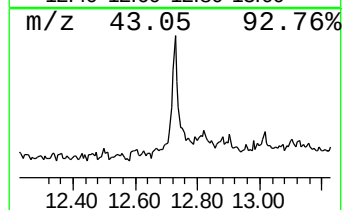
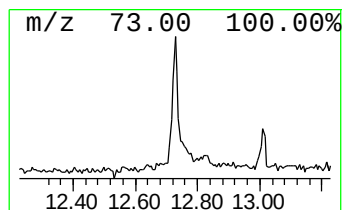
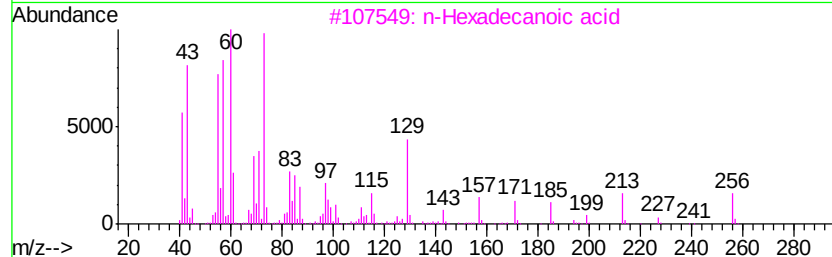
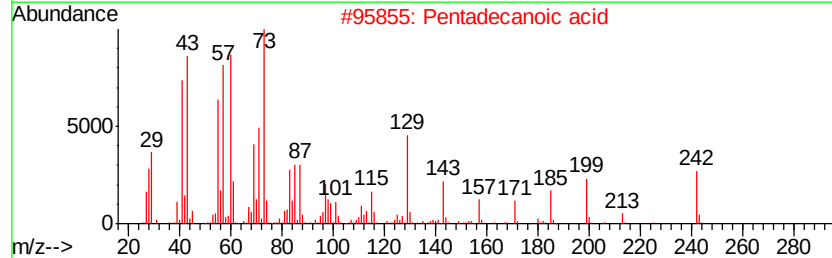
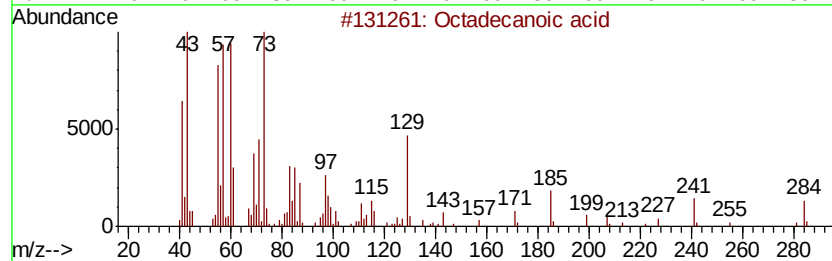
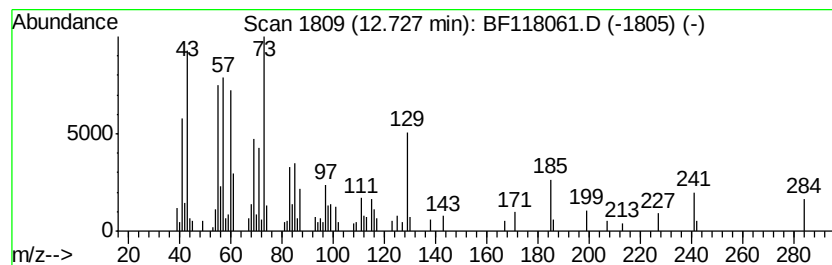
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Octadecanoic acid Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.73 | 2.17 ng | 143864 | Phenanthrene-d10 | 11.42 |

| Hit# of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|---------|---|---------------------|-----|----------|-------------|------|
| 1       |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 98   |
| 2       |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 89   |
| 3       |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 52   |
| 4       |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 46   |
| 5       |   | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 43   |



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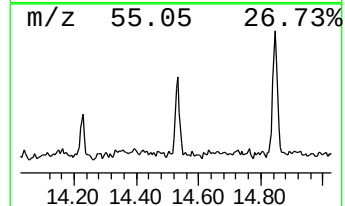
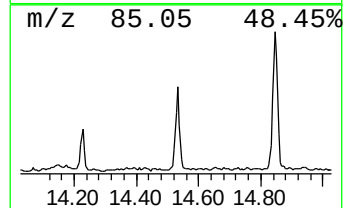
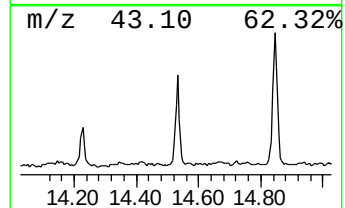
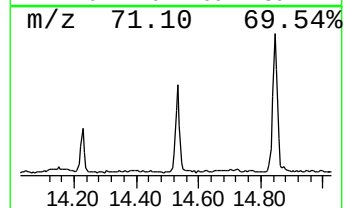
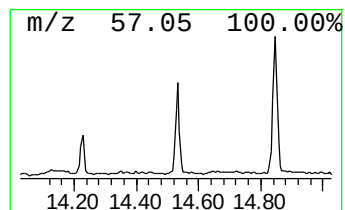
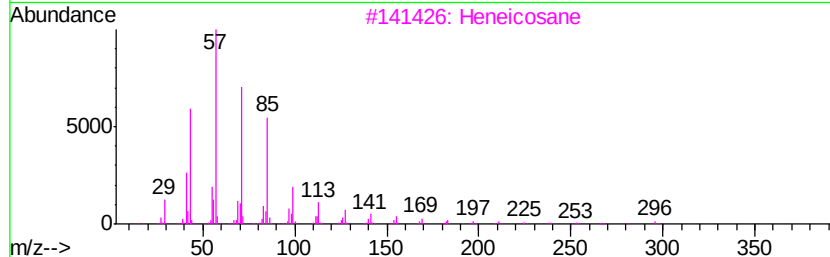
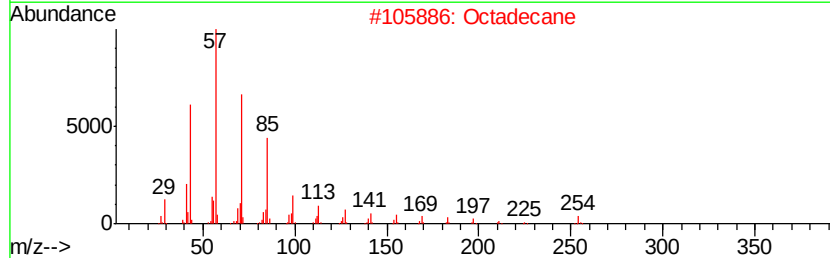
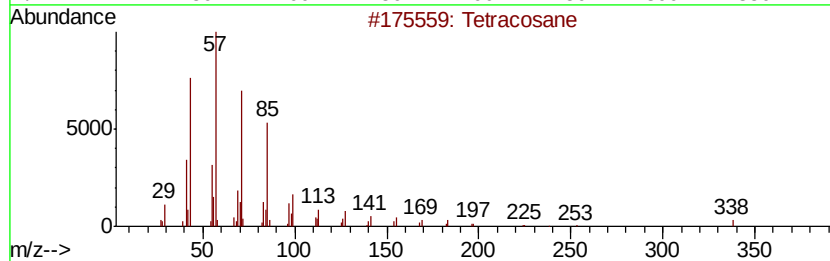
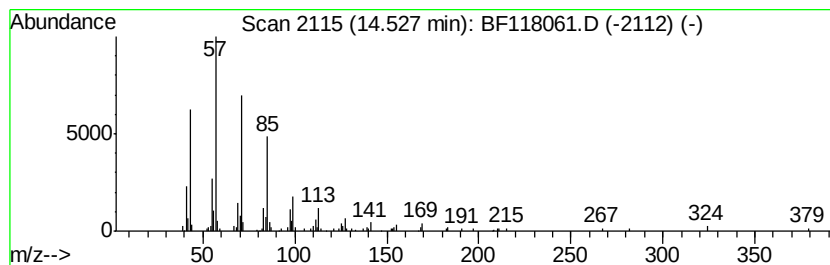
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Tetracosane Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.53 | 3.62 ng | 190166 | Chrysene-d12     | 14.07 |

| Hit# of 5 | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|-----------|-------------------|-----|---------|-------------|------|
| 1         | Tetracosane       | 338 | C24H50  | 000646-31-1 | 91   |
| 2         | Octadecane        | 254 | C18H38  | 000593-45-3 | 91   |
| 3         | Heneicosane       | 296 | C21H44  | 000629-94-7 | 91   |
| 4         | Triacontane       | 422 | C30H62  | 000638-68-6 | 91   |
| 5         | Tetratetracontane | 619 | C44H90  | 007098-22-8 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\  
Data File : BF118061.D  
Acq On : 3 Dec 2019 14:37  
Operator : JU  
Sample : PB125174BL  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB125174BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

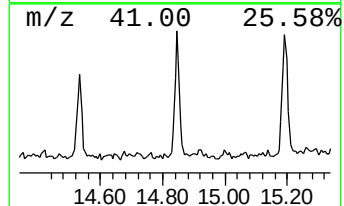
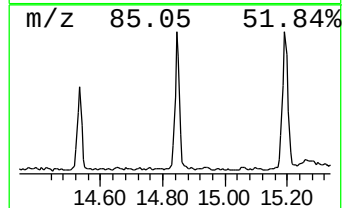
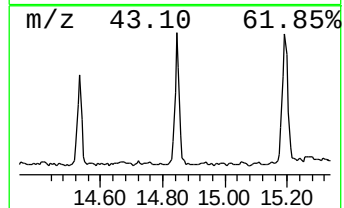
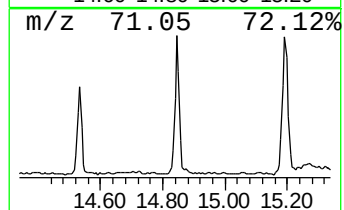
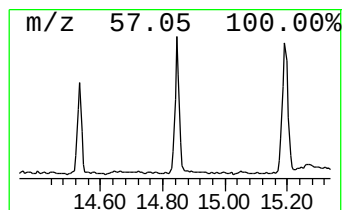
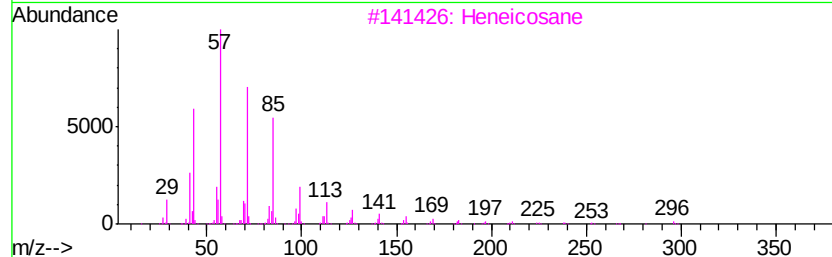
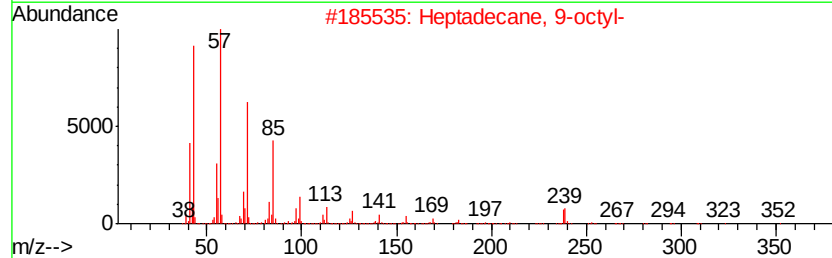
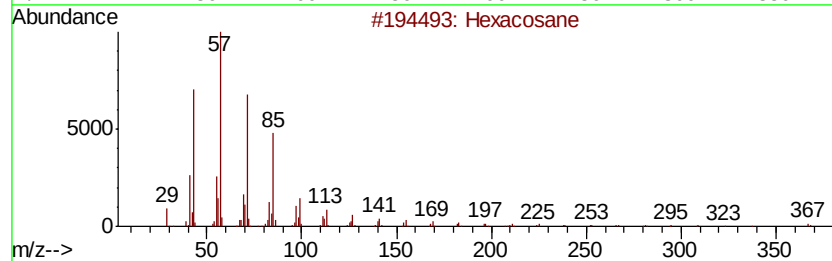
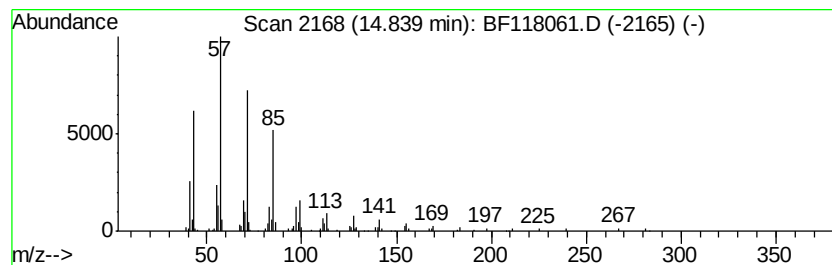
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Hexacosane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.84 | 4.70 ng | 308875 | Perylene-d12     | 15.56 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexacosane                       | 366 | C26H54  | 000630-01-3 | 94   |
| 2    |    | Heptadecane, 9-octyl-            | 352 | C25H52  | 007225-64-1 | 91   |
| 3    |    | Heneicosane                      | 296 | C21H44  | 000629-94-7 | 91   |
| 4    |    | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 91   |
| 5    |    | Docosane                         | 310 | C22H46  | 000629-97-0 | 91   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\  
Data File : BF118061.D  
Acq On : 3 Dec 2019 14:37  
Operator : JU  
Sample : PB125174BL  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB125174BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

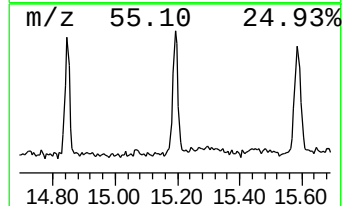
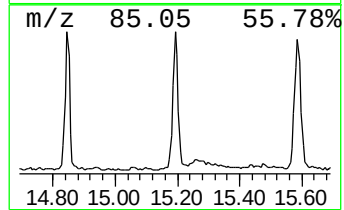
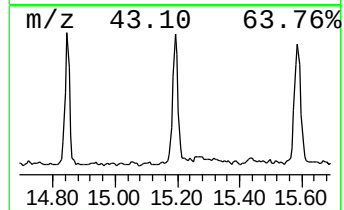
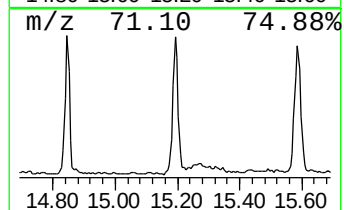
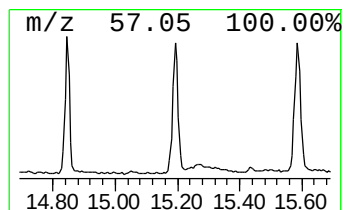
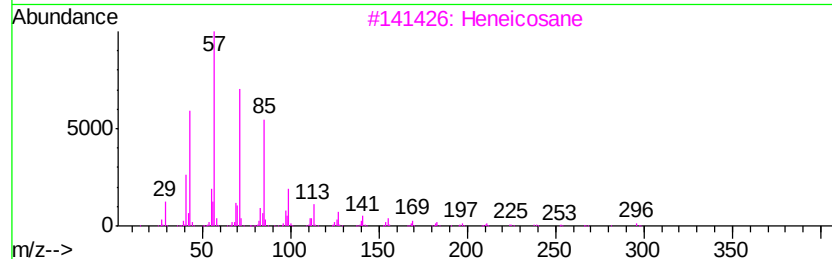
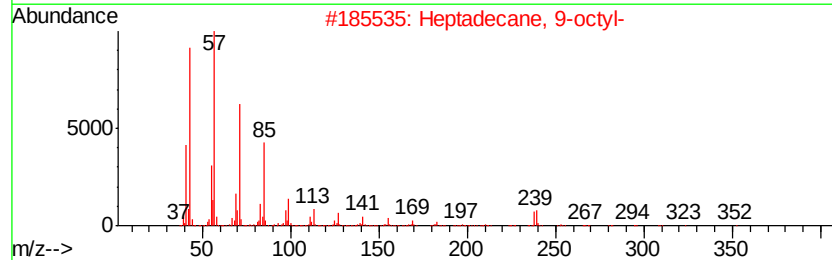
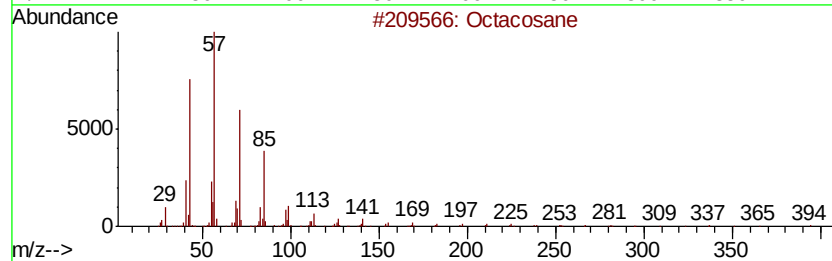
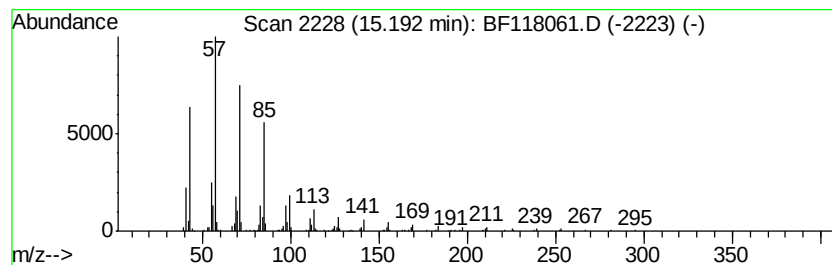
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Octacosane Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.19 | 5.92 ng | 389167 | Perylene-d12     | 15.56 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Octacosane            | 394 | C28H58  | 000630-02-4 | 91   |
| 2    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |
| 3    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 4    |    | Heptacosane           | 380 | C27H56  | 000593-49-7 | 91   |
| 5    |    | Heptadecane           | 240 | C17H36  | 000629-78-7 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\  
Data File : BF118061.D  
Acq On : 3 Dec 2019 14:37  
Operator : JU  
Sample : PB125174BL  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB125174BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

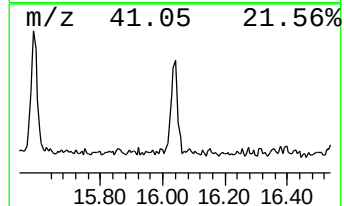
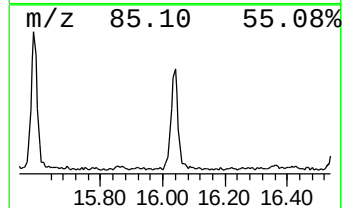
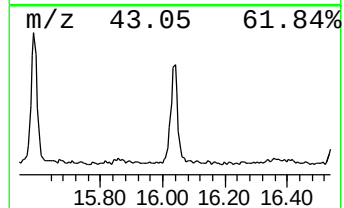
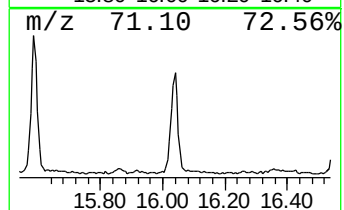
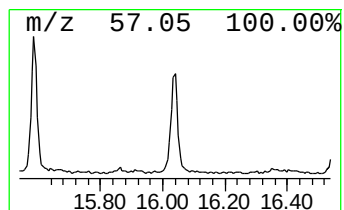
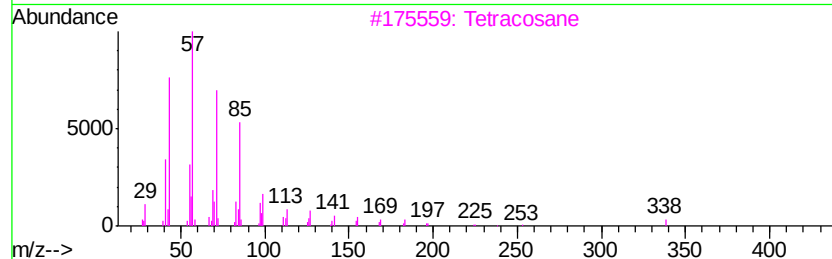
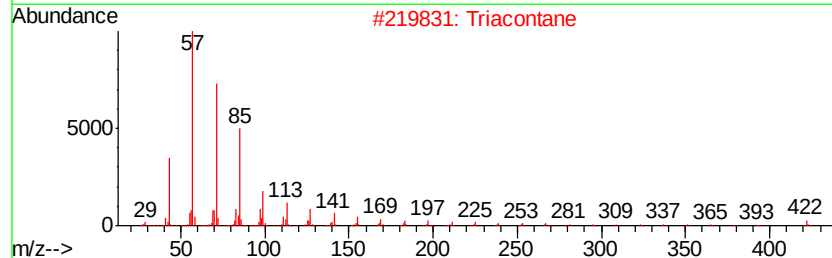
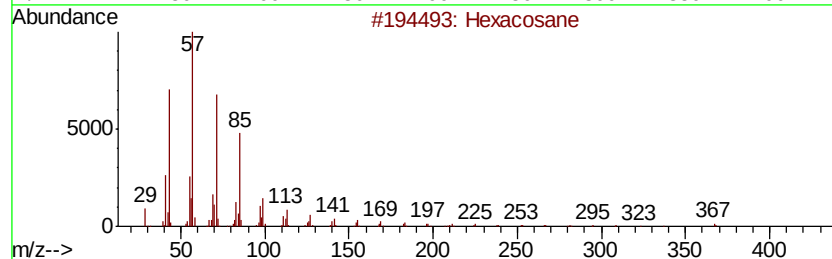
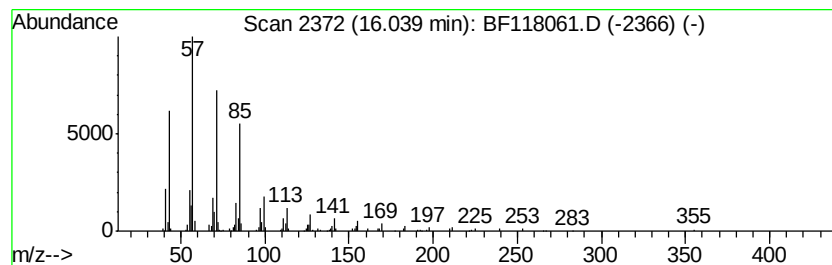
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Triacontane Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.04 | 5.08 ng | 334101 | Perylene-d12     | 15.56 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Hexacosane            | 366 | C26H54  | 000630-01-3 | 93   |
| 2    |    | Triacontane           | 422 | C30H62  | 000638-68-6 | 91   |
| 3    |    | Tetracosane           | 338 | C24H50  | 000646-31-1 | 91   |
| 4    |    | Tetratetracontane     | 619 | C44H90  | 007098-22-8 | 91   |
| 5    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120319\  
Data File : BF118061.D  
Acq On : 3 Dec 2019 14:37  
Operator : JU  
Sample : PB125174BL  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB125174BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

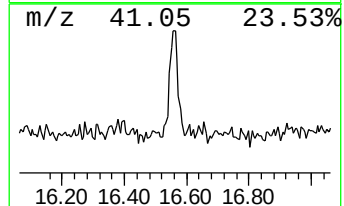
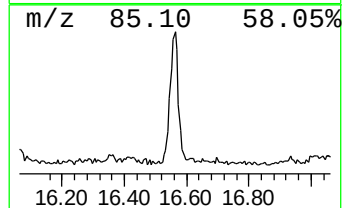
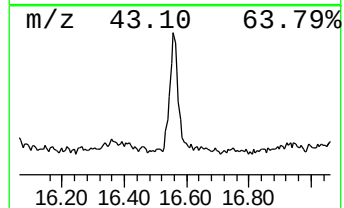
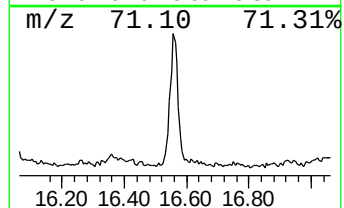
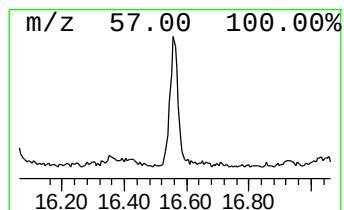
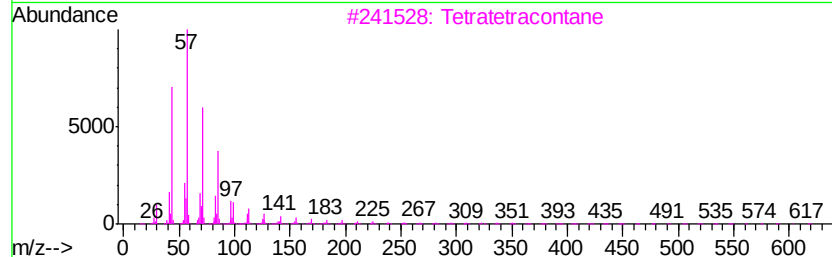
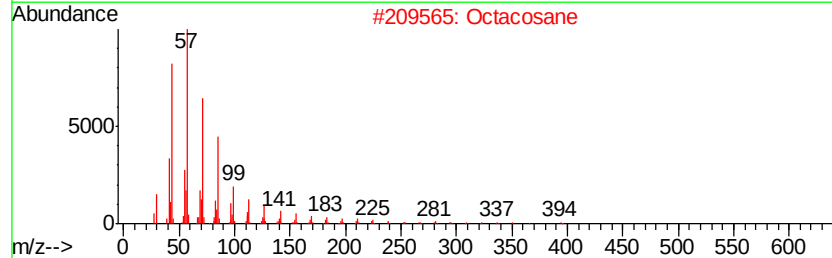
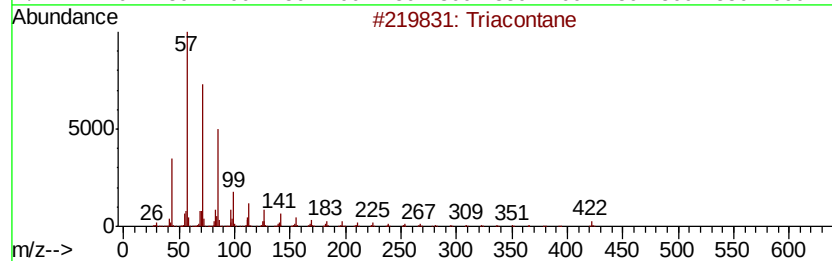
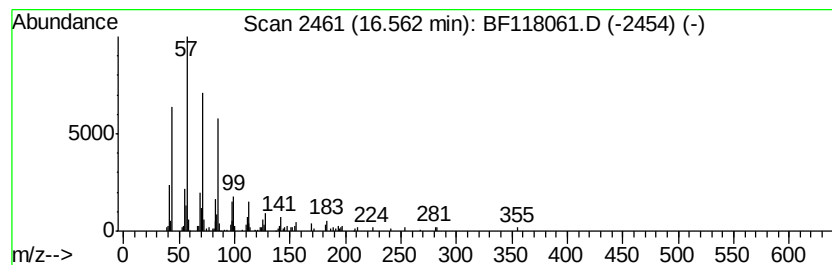
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Tetratetracontane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.56 | 3.69 ng | 242902 | Perylene-d12     | 15.56 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Triacontane       | 422 | C30H62  | 000638-68-6 | 91   |
| 2    |    | Octacosane        | 394 | C28H58  | 000630-02-4 | 91   |
| 3    |    | Tetratetracontane | 619 | C44H90  | 007098-22-8 | 91   |
| 4    |    | Heptacosane       | 380 | C27H56  | 000593-49-7 | 91   |
| 5    |    | Docosane          | 310 | C22H46  | 000629-97-0 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF120319\  
Data File : BF118061.D  
Acq On : 3 Dec 2019 14:37  
Operator : JU  
Sample : PB125174BL  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PB125174BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF111319.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... | 2.15  | 121.0   | ng    | 5005520  | 1                     | 6.87  | 827700  | 20.0 |
| 2-Pentanone, 4-hy... | 5.09  | 8.3     | ng    | 343702   | 1                     | 6.87  | 827700  | 20.0 |
| unknown6.65          | 6.65  | 96.8    | ng    | 4005260  | 1                     | 6.87  | 827700  | 20.0 |
| Octadecanoic acid    | 12.73 | 2.2     | ng    | 143864   | 4                     | 11.42 | 1328960 | 20.0 |
| Tetracosane          | 14.53 | 3.6     | ng    | 190166   | 5                     | 14.07 | 1050490 | 20.0 |
| Hexacosane           | 14.84 | 4.7     | ng    | 308875   | 6                     | 15.56 | 1315350 | 20.0 |
| Octacosane           | 15.19 | 5.9     | ng    | 389167   | 6                     | 15.56 | 1315350 | 20.0 |
| Triacontane          | 16.04 | 5.1     | ng    | 334101   | 6                     | 15.56 | 1315350 | 20.0 |
| Tetratetracontane    | 16.56 | 3.7     | ng    | 242902   | 6                     | 15.56 | 1315350 | 20.0 |