

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121119\  
 Data File : BP001315.D  
 Acq On : 11 Dec 2019 16:59  
 Operator : JU  
 Sample : K6204-07  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 K072-AA-WC-2

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\_P\METHODS\8270-BP112719.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 5.616       | 591           | 600         | 612          | rBV      | 229074         | 390334        | 11.67%          | 1.811%        |
| 2         | 6.216       | 695           | 702         | 718          | rBV      | 1447148        | 1850467       | 55.33%          | 8.587%        |
| 3         | 7.757       | 957           | 964         | 978          | rBV      | 1560762        | 2107192       | 63.01%          | 9.778%        |
| 4         | 7.945       | 989           | 996         | 1011         | rBV      | 1676802        | 2013242       | 60.20%          | 9.342%        |
| 5         | 8.304       | 1052          | 1057        | 1064         | rVB      | 391444         | 427071        | 12.77%          | 1.982%        |
| 6         | 8.545       | 1092          | 1098        | 1103         | rBV      | 1318744        | 1528982       | 45.72%          | 7.095%        |
| 7         | 9.187       | 1200          | 1207        | 1211         | rBV      | 1078061        | 1301851       | 38.93%          | 6.041%        |
| 8         | 10.339      | 1398          | 1403        | 1414         | rBV      | 459155         | 526546        | 15.74%          | 2.443%        |
| 9         | 12.122      | 1699          | 1706        | 1710         | rBV      | 2216928        | 2506807       | 74.95%          | 11.632%       |
| 10        | 12.786      | 1814          | 1819        | 1826         | rBV      | 80041          | 96235         | 2.88%           | 0.447%        |
| 11        | 13.204      | 1884          | 1890        | 1905         | rBV      | 526906         | 673334        | 20.13%          | 3.124%        |
| 12        | 14.498      | 2103          | 2110        | 2133         | rBV      | 1509581        | 1925466       | 57.57%          | 8.935%        |
| 13        | 15.598      | 2288          | 2297        | 2309         | rBV      | 648977         | 790454        | 23.63%          | 3.668%        |
| 14        | 16.492      | 2445          | 2449        | 2464         | rBV2     | 108218         | 161316        | 4.82%           | 0.749%        |
| 15        | 17.574      | 2628          | 2633        | 2643         | rBV2     | 21666          | 37462         | 1.12%           | 0.174%        |
| 16        | 17.886      | 2680          | 2686        | 2690         | rBV      | 3049426        | 3344480       | 100.00%         | 15.519%       |
| 17        | 19.157      | 2892          | 2902        | 2911         | rBV7     | 61218          | 263979        | 7.89%           | 1.225%        |
| 18        | 19.239      | 2911          | 2916        | 2930         | rVB      | 639414         | 772948        | 23.11%          | 3.587%        |
| 19        | 19.921      | 3028          | 3032        | 3037         | rBV      | 38769          | 39367         | 1.18%           | 0.183%        |
| 20        | 20.298      | 3092          | 3096        | 3100         | rBV      | 44151          | 45632         | 1.36%           | 0.212%        |
| 21        | 20.692      | 3158          | 3163        | 3167         | rBV      | 43978          | 56265         | 1.68%           | 0.261%        |
| 22        | 21.009      | 3211          | 3217        | 3227         | rBV      | 409221         | 577607        | 17.27%          | 2.680%        |
| 23        | 21.127      | 3232          | 3237        | 3243         | rBV2     | 37786          | 56013         | 1.67%           | 0.260%        |
| 24        | 21.621      | 3315          | 3321        | 3327         | rBV2     | 34225          | 57526         | 1.72%           | 0.267%        |

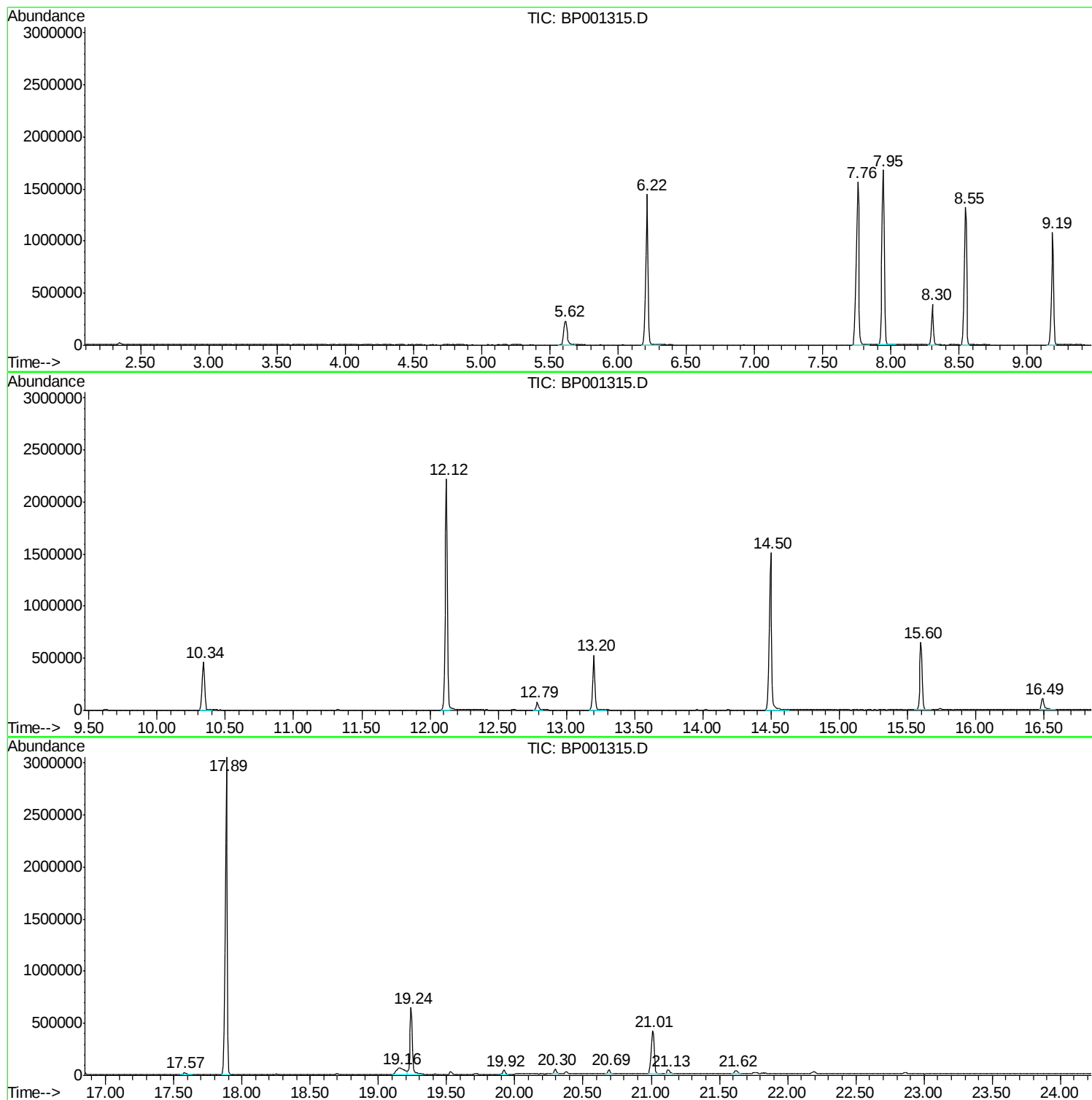
Sum of corrected areas: 21550576

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121119\  
Data File : BP001315.D  
Acq On : 11 Dec 2019 16:59  
Operator : JU  
Sample : K6204-07  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
K072-AA-WC-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121119\  
Data File : BP001315.D  
Acq On : 11 Dec 2019 16:59  
Operator : JU  
Sample : K6204-07  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
K072-AA-WC-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

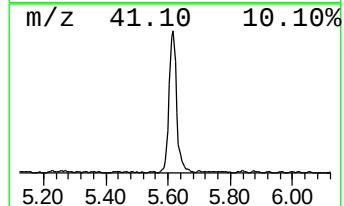
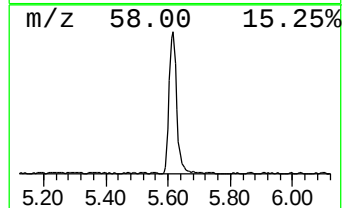
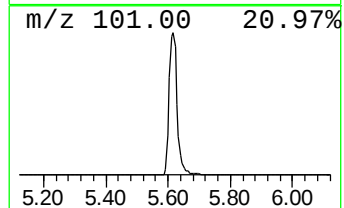
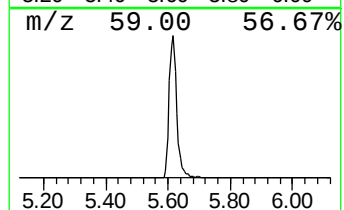
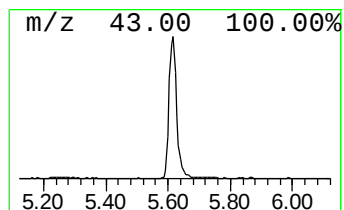
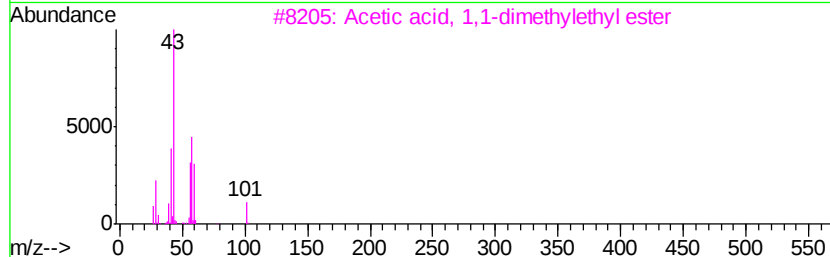
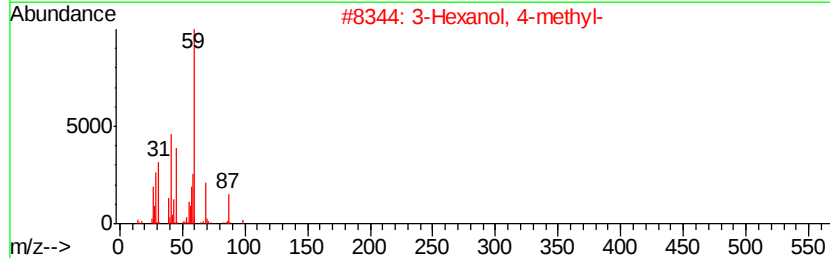
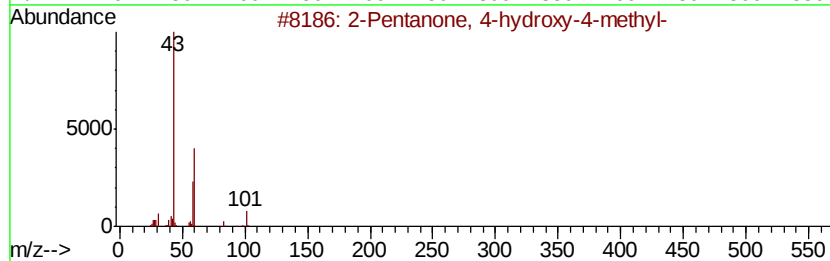
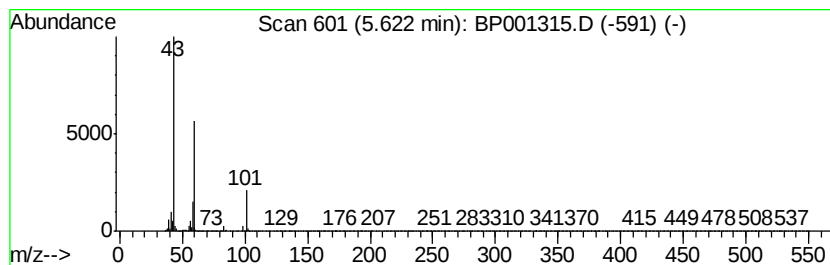
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.62 | 18.28 ng | 390334 | 1,4-Dichlorobenzene-d4 | 8.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2 | 45   |
| 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    |    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2 | 000057-71-6 | 27   |
| 5    |    | 2-Hexanol, 2-methyl-                | 116 | C7H16O  | 000625-23-0 | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121119\  
Data File : BP001315.D  
Acq On : 11 Dec 2019 16:59  
Operator : JU  
Sample : K6204-07  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
K072-AA-WC-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

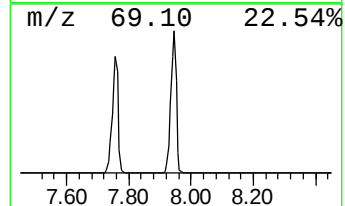
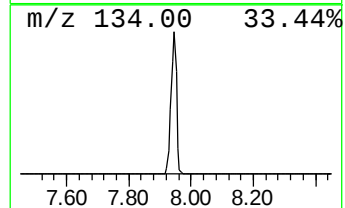
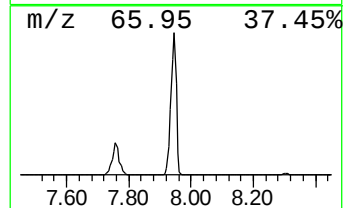
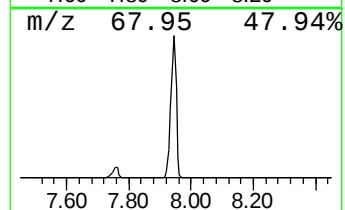
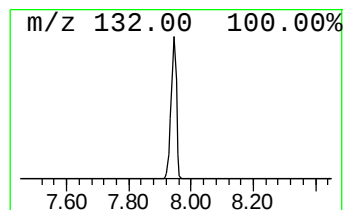
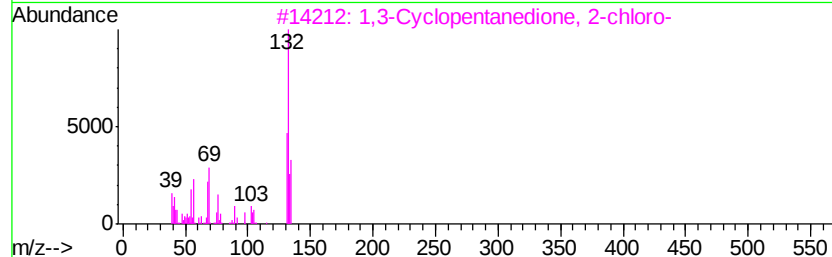
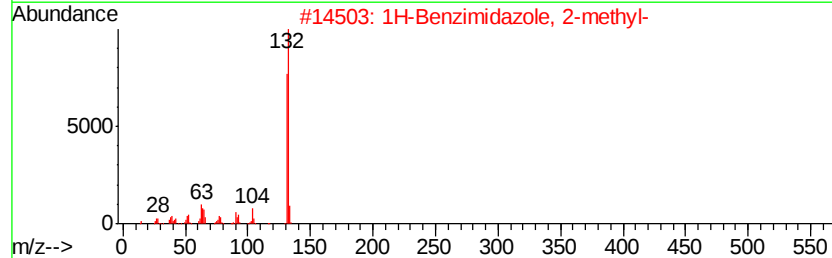
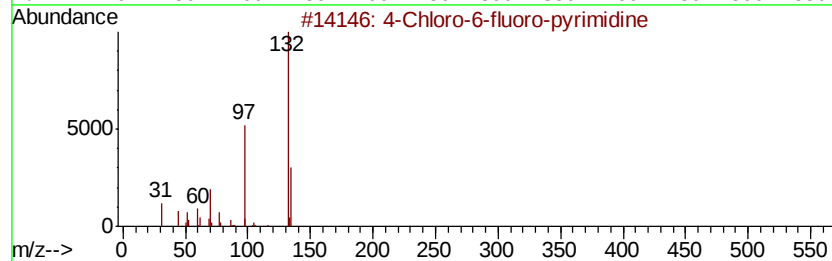
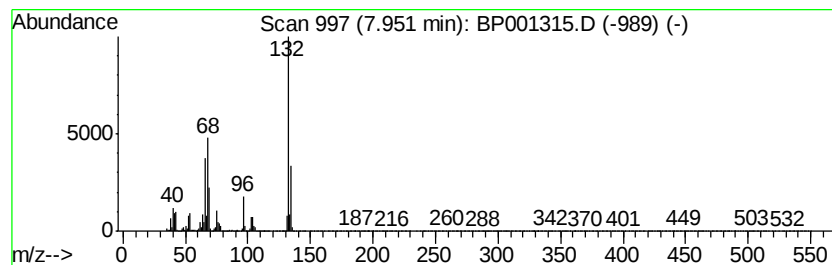
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown7.95 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.95 | 94.28 ng | 2013240 | 1,4-Dichlorobenzene-d4 | 8.30 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4-Chloro-6-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-01-6  | 14   |
| 2    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 14   |
| 3    |    | 1,3-Cyclopentanedione, 2-chloro- | 132 | C5H5ClO2  | 014203-19-1  | 12   |
| 4    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 10   |
| 5    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121119\  
Data File : BP001315.D  
Acq On : 11 Dec 2019 16:59  
Operator : JU  
Sample : K6204-07  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

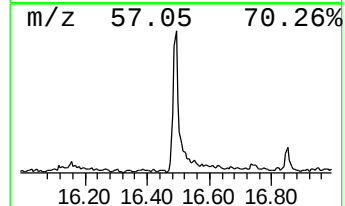
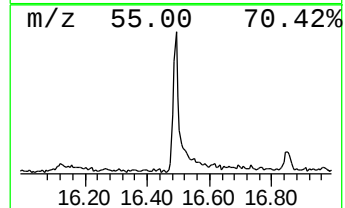
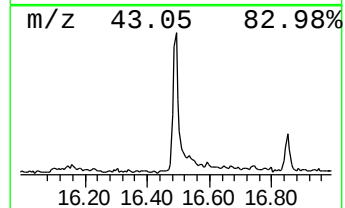
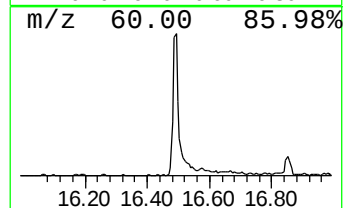
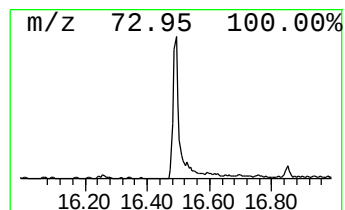
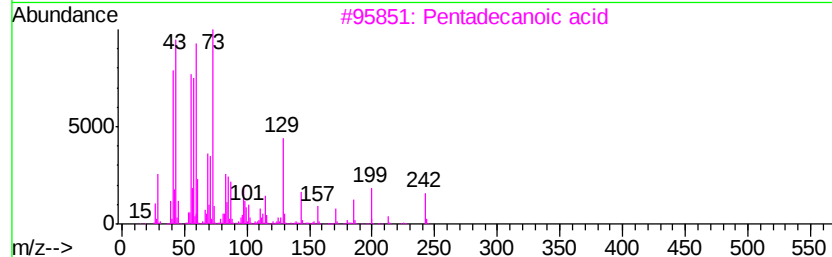
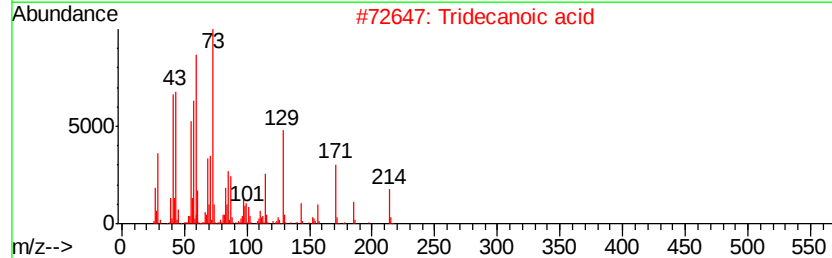
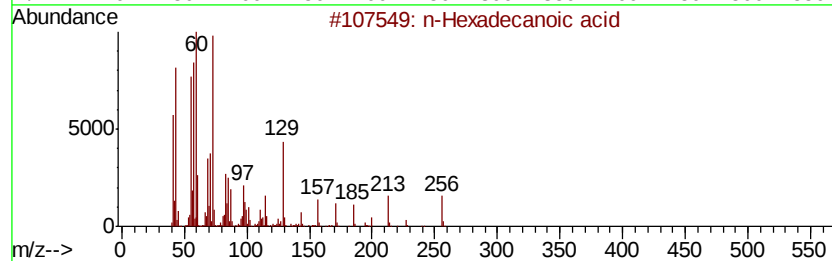
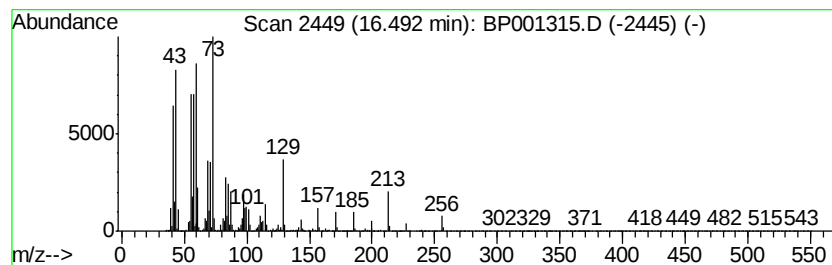
Instrument :  
BNA\_P  
ClientSampleId :  
K072-AA-WC-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

| R.T.    | EstConc             | Area         | Relative to ISTD |          | R.T.        |      |
|---------|---------------------|--------------|------------------|----------|-------------|------|
| 16.49   | 4.08 ng             | 161316       | Phenanthrene-d10 |          | 15.60       |      |
| Hit# of | 5                   | Tentative ID | MW               | MolForm  | CAS#        | Qual |
| 1       | n-Hexadecanoic acid |              | 256              | C16H32O2 | 000057-10-3 | 97   |
| 2       | Tridecanoic acid    |              | 214              | C13H26O2 | 000638-53-9 | 90   |
| 3       | Pentadecanoic acid  |              | 242              | C15H30O2 | 001002-84-2 | 87   |
| 4       | Tetradecanoic acid  |              | 228              | C14H28O2 | 000544-63-8 | 81   |
| 5       | Dodecanoic acid     |              | 200              | C12H24O2 | 000143-07-7 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121119\  
Data File : BP001315.D  
Acq On : 11 Dec 2019 16:59  
Operator : JU  
Sample : K6204-07  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
K072-AA-WC-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

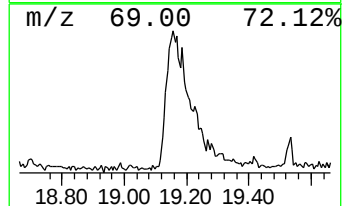
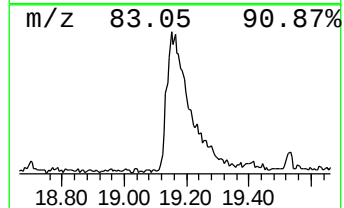
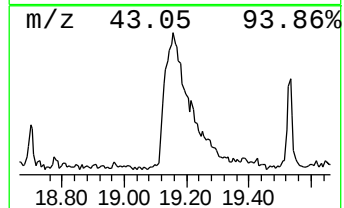
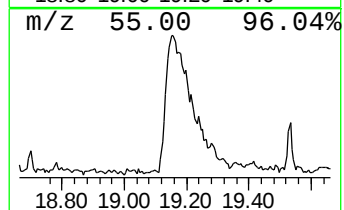
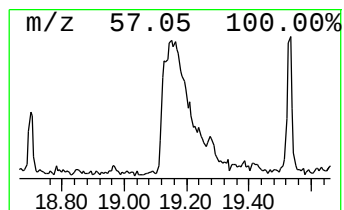
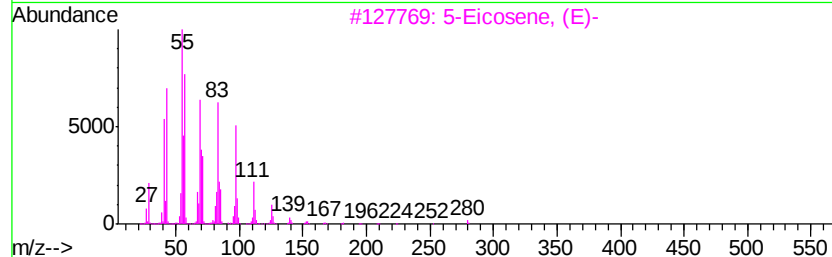
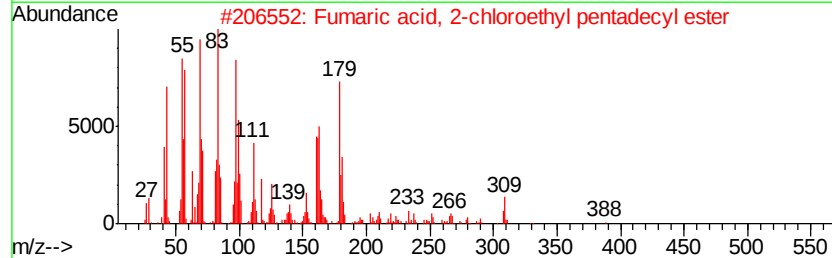
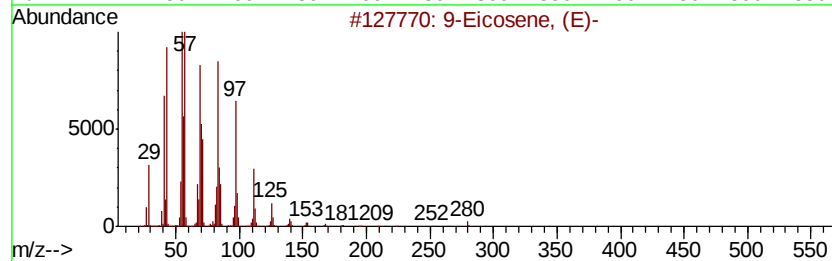
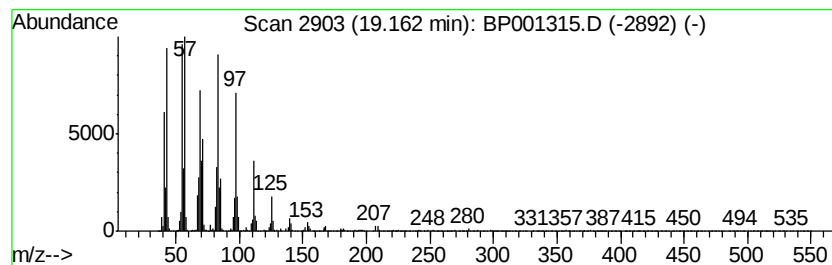
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 9-Eicosene, (E)- Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.16 | 6.83 ng | 263979 | Chrysene-d12     | 19.24 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm    | CAS#         | Qual |
|---------|-------------------------------------|--------------|-----|------------|--------------|------|
| 1       | 9-Eicosene, (E)-                    |              | 280 | C20H40     | 074685-29-3  | 96   |
| 2       | Fumaric acid, 2-chloroethyl pent... |              | 388 | C21H37ClO4 | 1000330-62-1 | 96   |
| 3       | 5-Eicosene, (E)-                    |              | 280 | C20H40     | 074685-30-6  | 96   |
| 4       | Fumaric acid, 2-chloroethyl dode... |              | 346 | C18H31ClO4 | 1000330-62-0 | 96   |
| 5       | 9-Tricosene, (Z)-                   |              | 322 | C23H46     | 027519-02-4  | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP121119\  
Data File : BP001315.D  
Acq On : 11 Dec 2019 16:59  
Operator : JU  
Sample : K6204-07  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
K072-AA-WC-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP112719.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... | 5.62  | 18.3    | ng    | 390334   | 1                     | 8.30  | 427071 | 20.0 |
| unknown7.95          | 7.95  | 94.3    | ng    | 2013240  | 1                     | 8.30  | 427071 | 20.0 |
| n-Hexadecanoic acid  | 16.49 | 4.1     | ng    | 161316   | 4                     | 15.60 | 790454 | 20.0 |
| 9-Eicosene, (E)-     | 19.16 | 6.8     | ng    | 263979   | 5                     | 19.24 | 772948 | 20.0 |