

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
 Data File : BG023117.D  
 Acq On : 15 Jul 2016 17:28  
 Operator : UM/SJ  
 Sample : H4017-17  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C5-9

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:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.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         | 3.040       | 30            | 34          | 44           | rBV      | 1083710        | 1502993       | 19.14%          | 3.113%        |
| 2         | 5.226       | 399           | 406         | 411          | rBV      | 228346         | 363761        | 4.63%           | 0.753%        |
| 3         | 5.696       | 480           | 486         | 500          | rBV      | 1568320        | 2722133       | 34.66%          | 5.637%        |
| 4         | 7.306       | 753           | 760         | 774          | rVB      | 1845832        | 3363838       | 42.83%          | 6.966%        |
| 5         | 7.682       | 817           | 824         | 836          | rBV      | 1740024        | 3095091       | 39.41%          | 6.410%        |
| 6         | 8.152       | 896           | 904         | 911          | rVB      | 417819         | 722471        | 9.20%           | 1.496%        |
| 7         | 8.481       | 952           | 960         | 970          | rVB      | 1219137        | 2217829       | 28.24%          | 4.593%        |
| 8         | 9.327       | 1095          | 1104        | 1111         | rBV      | 1212507        | 2191282       | 27.90%          | 4.538%        |
| 9         | 10.978      | 1377          | 1385        | 1393         | rBV      | 590237         | 1102717       | 14.04%          | 2.284%        |
| 10        | 13.417      | 1793          | 1800        | 1812         | rBV      | 3210407        | 5219798       | 66.47%          | 10.810%       |
| 11        | 13.675      | 1839          | 1844        | 1849         | rVB4     | 52231          | 85988         | 1.09%           | 0.178%        |
| 12        | 14.239      | 1932          | 1940        | 1950         | rBV      | 1381317        | 2030960       | 25.86%          | 4.206%        |
| 13        | 14.803      | 2028          | 2036        | 2045         | rVB2     | 1044796        | 1716631       | 21.86%          | 3.555%        |
| 14        | 16.290      | 2282          | 2289        | 2301         | rBV      | 2732203        | 4428013       | 56.38%          | 9.170%        |
| 15        | 16.483      | 2318          | 2322        | 2324         | rBV4     | 72331          | 88283         | 1.12%           | 0.183%        |
| 16        | 17.553      | 2496          | 2504        | 2513         | rBV2     | 1431375        | 2200688       | 28.02%          | 4.557%        |
| 17        | 17.911      | 2559          | 2565        | 2575         | rBV3     | 180380         | 430238        | 5.48%           | 0.891%        |
| 18        | 18.417      | 2646          | 2651        | 2657         | rBV2     | 157038         | 222145        | 2.83%           | 0.460%        |
| 19        | 19.274      | 2792          | 2797        | 2815         | rBV      | 555515         | 1249794       | 15.91%          | 2.588%        |
| 20        | 20.150      | 2940          | 2946        | 2956         | rBV      | 5790130        | 7853271       | 100.00%         | 16.264%       |
| 21        | 20.773      | 3046          | 3052        | 3056         | rBV3     | 117787         | 208717        | 2.66%           | 0.432%        |
| 22        | 20.890      | 3069          | 3072        | 3076         | rVB3     | 69859          | 86266         | 1.10%           | 0.179%        |
| 23        | 21.472      | 3167          | 3171        | 3182         | rBV6     | 113815         | 309254        | 3.94%           | 0.640%        |
| 24        | 21.713      | 3208          | 3212        | 3218         | rVB2     | 79160          | 110886        | 1.41%           | 0.230%        |
| 25        | 21.854      | 3230          | 3236        | 3245         | rBV      | 1428609        | 2461578       | 31.34%          | 5.098%        |
| 26        | 25.232      | 3800          | 3811        | 3824         | rBV      | 711297         | 2302653       | 29.32%          | 4.769%        |

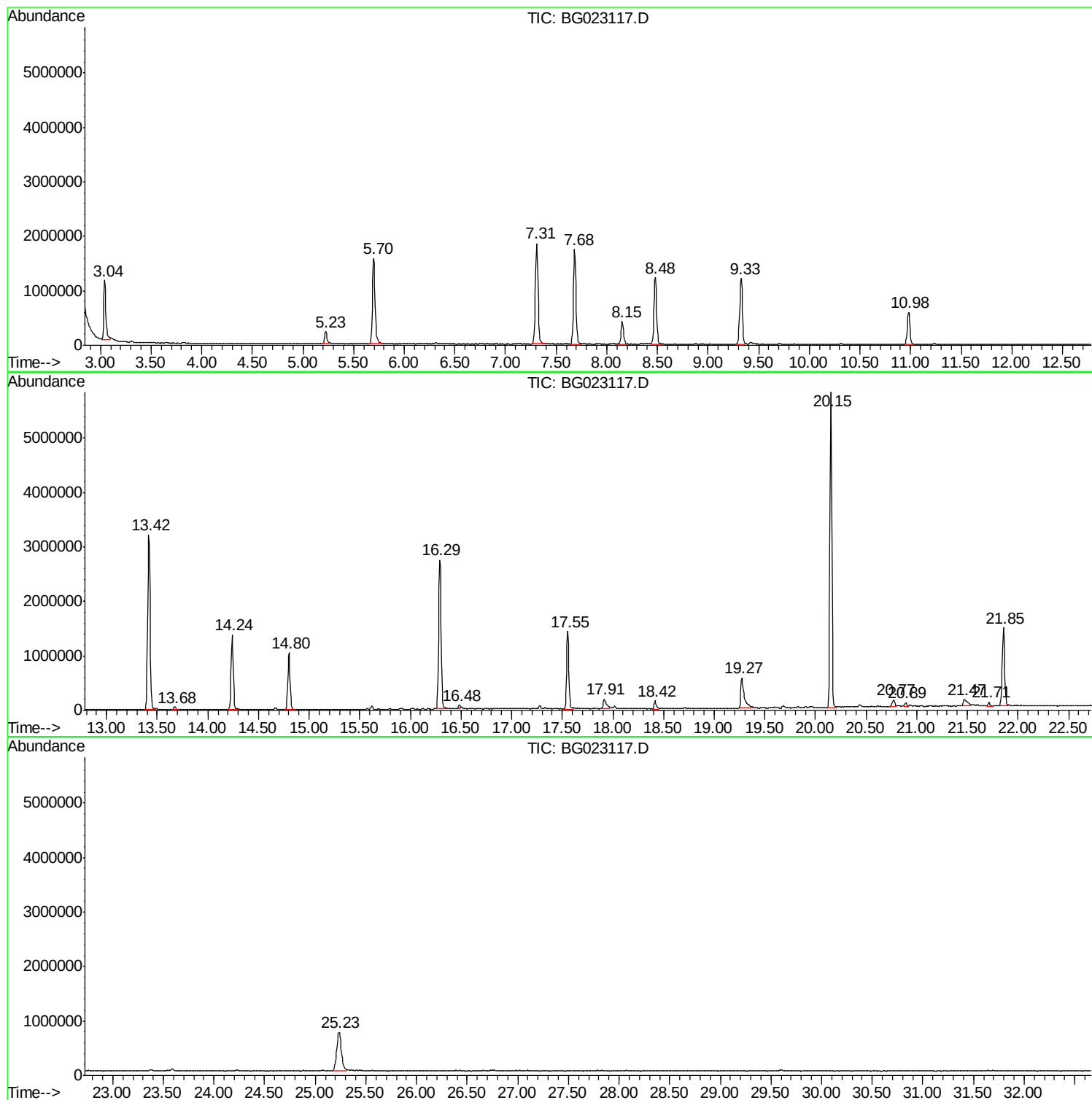
Sum of corrected areas: 48287278

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

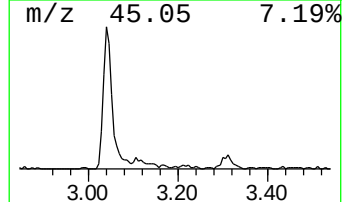
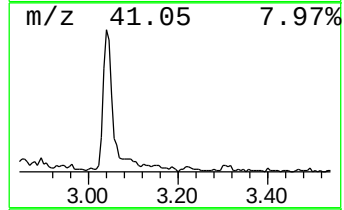
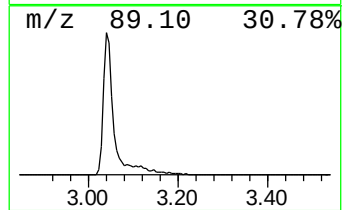
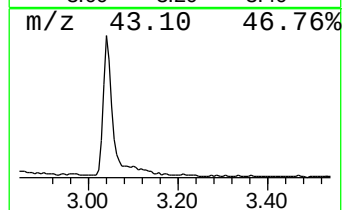
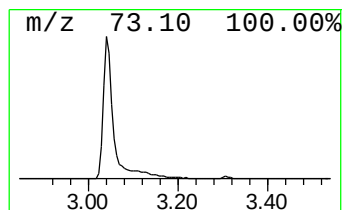
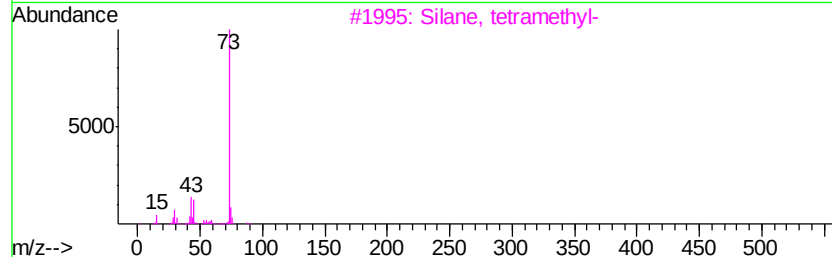
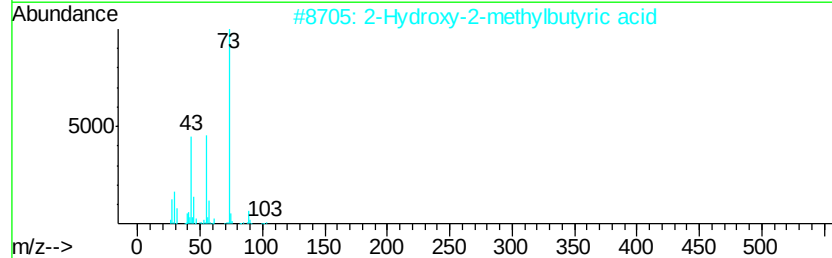
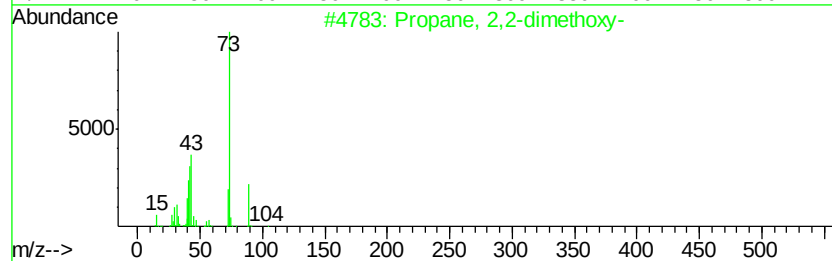
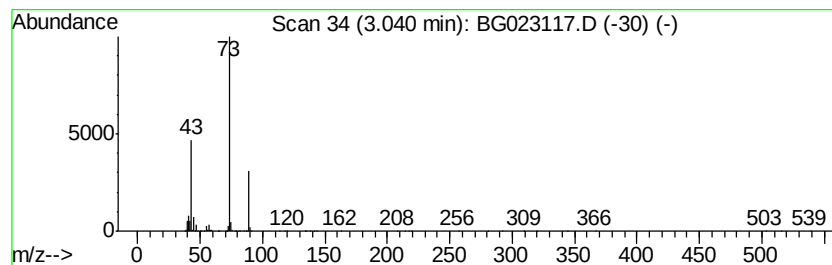
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 3.04 | 41.61 ng | 1502990 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Propane, 2,2-dimethoxy-             | 104 | C5H12O2  | 000077-76-9 | 72   |
| 2    |    |   | 2-Hydroxy-2-methylbutyric acid      | 118 | C5H10O3  | 003739-30-8 | 33   |
| 3    |    |   | Silane, tetramethyl-                | 88  | C4H12Si  | 000075-76-3 | 9    |
| 4    |    |   | 1-Hexen-4-ol, 1-chloro-3,5-dimet... | 162 | C8H15ClO | 100244-09-5 | 9    |
| 5    |    |   | Propane, 2-methoxy-2-methyl-        | 88  | C5H12O   | 001634-04-4 | 9    |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

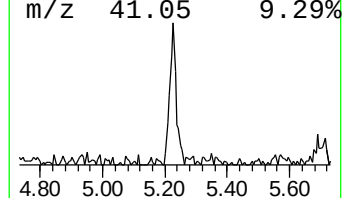
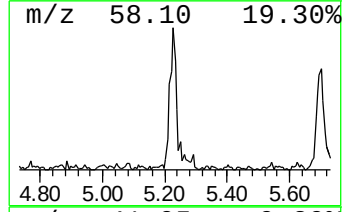
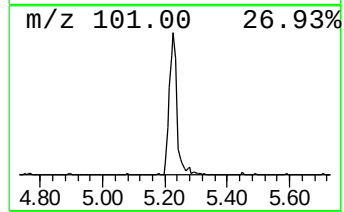
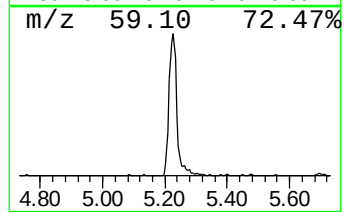
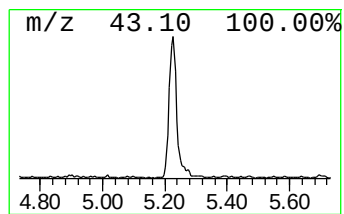
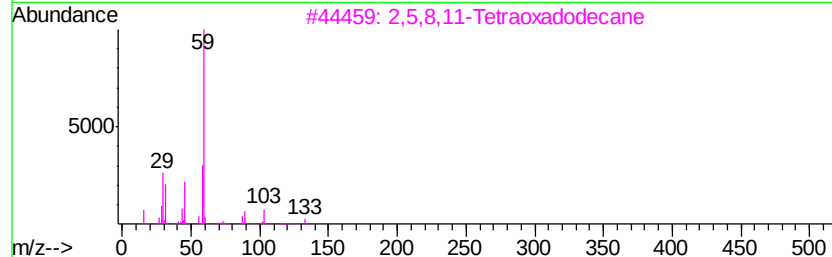
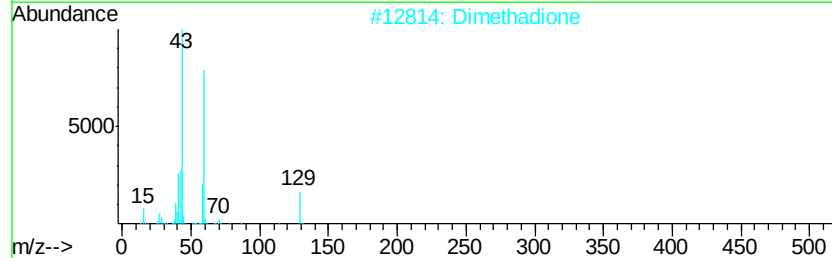
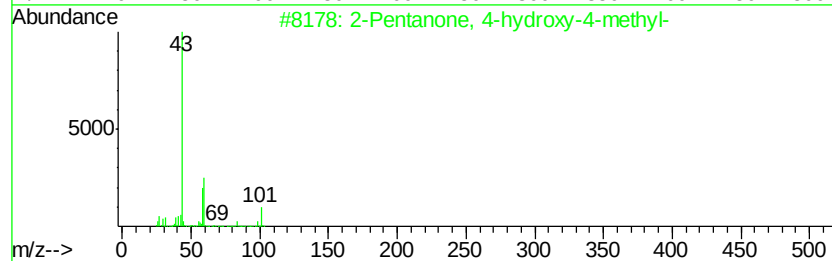
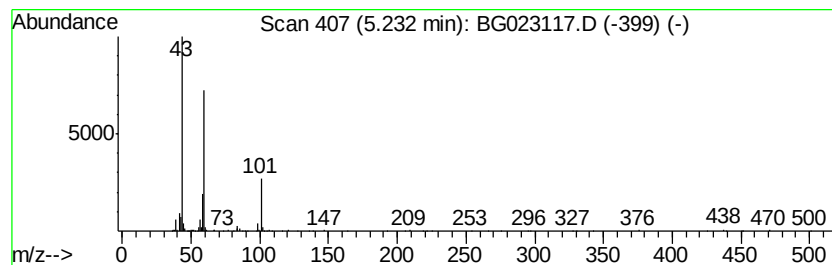
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown5.23 Concentration Rank 5

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.23 | 10.07 ng | 363761 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2  | 42   |
| 2    |    | Dimethadione                        | 129 | C5H7NO3  | 000695-53-4  | 36   |
| 3    |    | 2,5,8,11-Tetraoxadodecane           | 178 | C8H18O4  | 000112-49-2  | 25   |
| 4    |    | Succinic acid, heptyl 2-methoxye... | 274 | C14H26O5 | 1000325-80-7 | 25   |
| 5    |    | Silane, trimethyl-                  | 74  | C3H10Si  | 000993-07-7  | 23   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

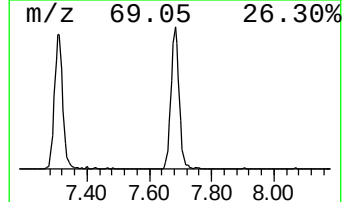
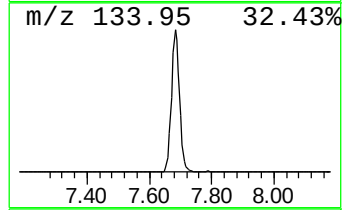
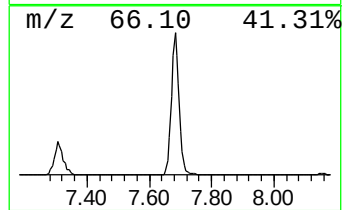
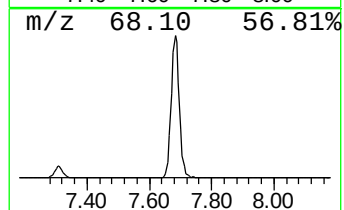
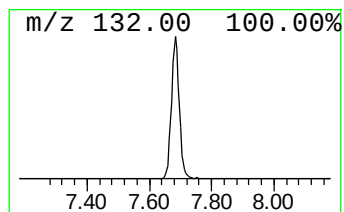
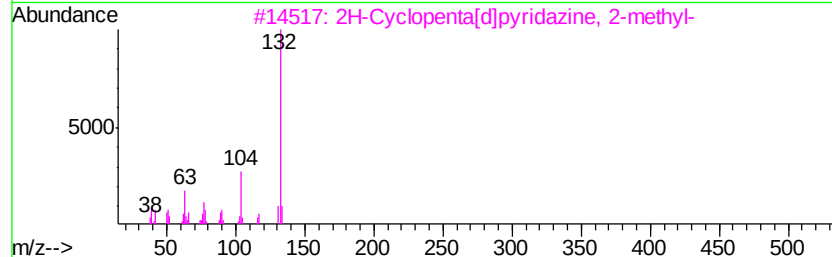
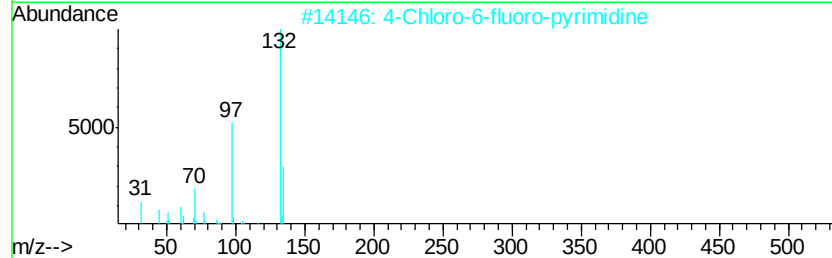
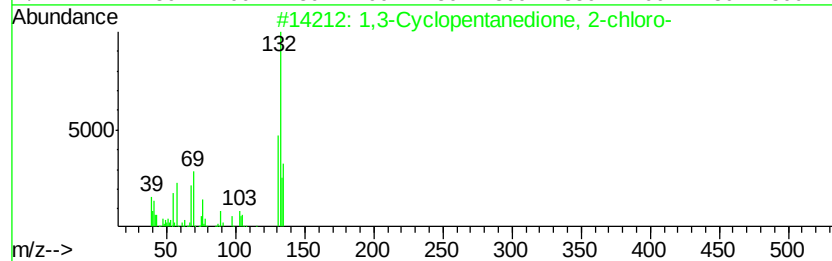
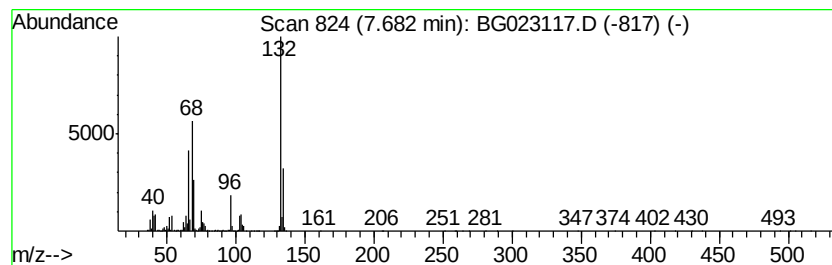
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown7.68 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.68 | 85.68 ng | 3095090 | 1,4-Dichlorobenzene-d4 | 8.15 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 1,3-Cyclopentanedione, 2-chloro-    | 132 | C5H5ClO2  | 014203-19-1  | 16   |
| 2    |    | 4-Chloro-6-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-01-6  | 14   |
| 3    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2    | 022291-85-6  | 14   |
| 4    |    | 1,3,5-Trifluorobenzene              | 132 | C6H3F3    | 000372-38-3  | 10   |
| 5    |    | Cyclopropanamine, 1-phenyl-         | 133 | C9H11N    | 1000351-26-2 | 10   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

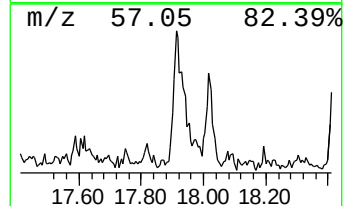
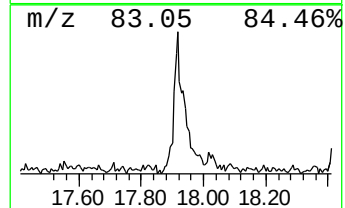
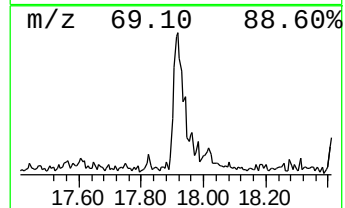
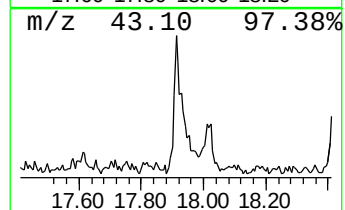
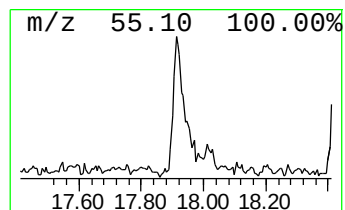
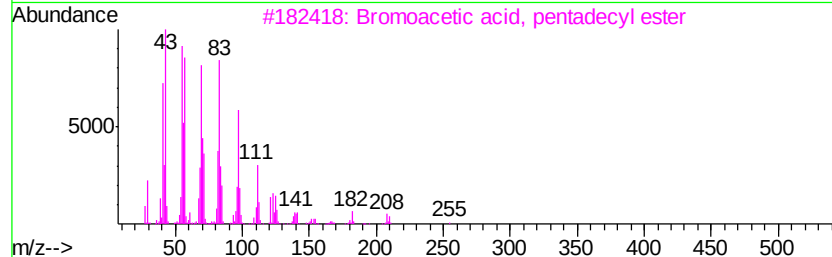
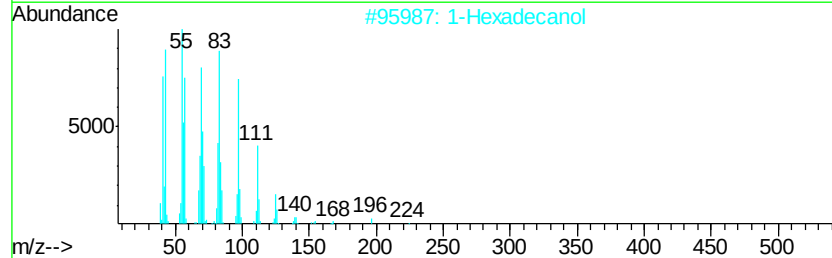
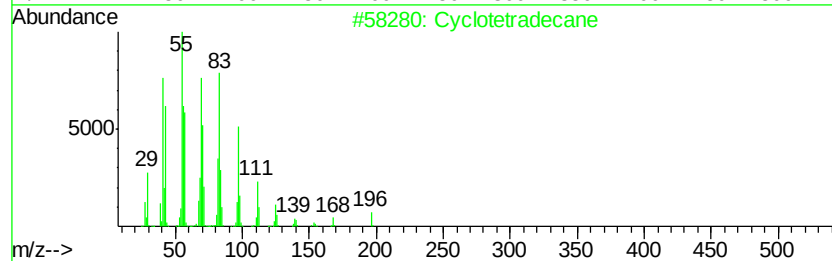
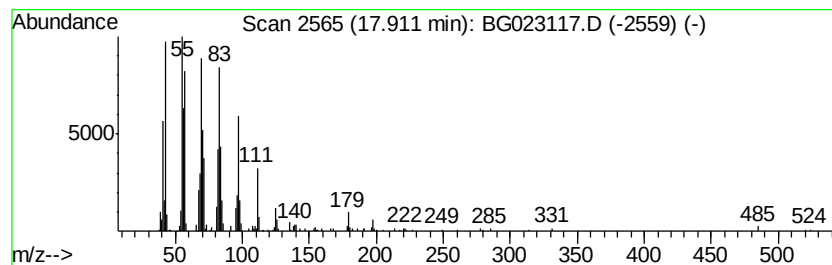
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Cyclotetradecane Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.91 | 3.91 ng | 430238 | Phenanthrene-d10 | 17.55 |

| Hit# of | 5 | Tentative ID                       | MW  | MolForm    | CAS#        | Qual |
|---------|---|------------------------------------|-----|------------|-------------|------|
| 1       |   | Cyclotetradecane                   | 196 | C14H28     | 000295-17-0 | 94   |
| 2       |   | 1-Hexadecanol                      | 242 | C16H34O    | 036653-82-4 | 90   |
| 3       |   | Bromoacetic acid, pentadecyl ester | 348 | C17H33BrO2 | 131143-01-6 | 90   |
| 4       |   | n-Heptadecanol-1                   | 256 | C17H36O    | 001454-85-9 | 87   |
| 5       |   | Cyclohexadecane                    | 224 | C16H32     | 000295-65-8 | 87   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

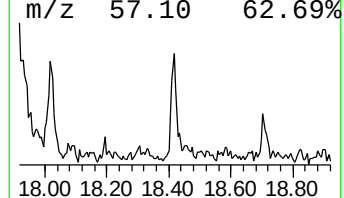
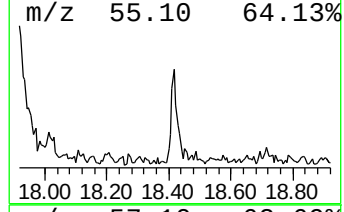
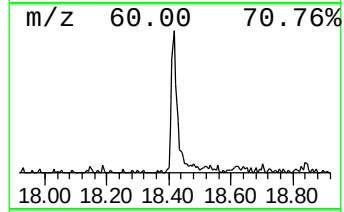
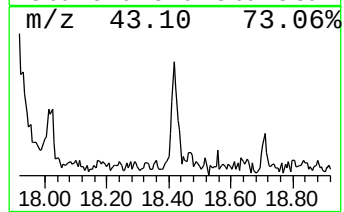
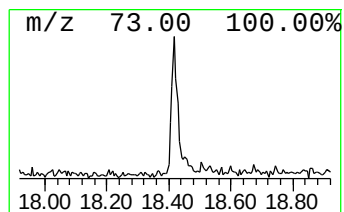
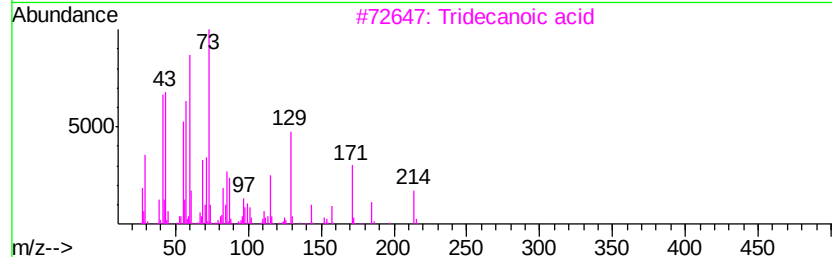
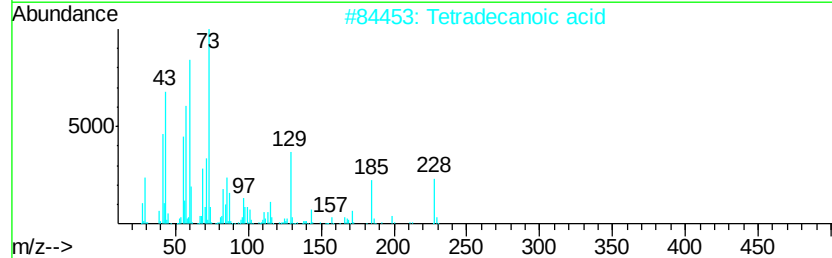
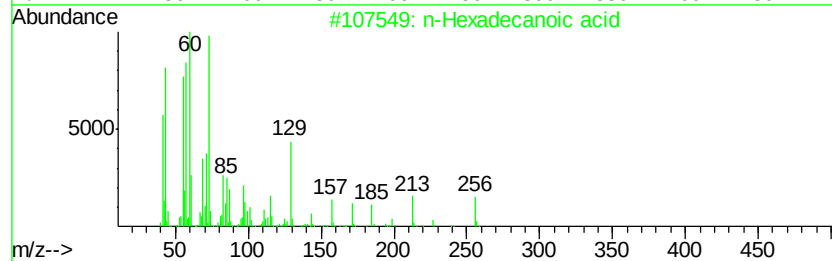
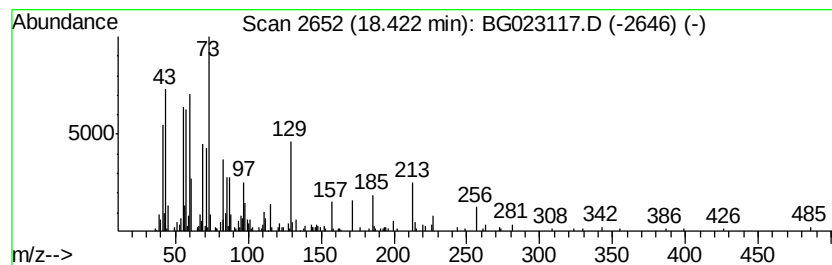
Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 n-Hexadecanoic acid Concentration Rank 8

| R.T.    | EstConc | Area                | Relative to ISTD |          | R.T.        |      |
|---------|---------|---------------------|------------------|----------|-------------|------|
| 18.42   | 2.02 ng | 222145              | Phenanthrene-d10 |          | 17.55       |      |
| Hit# of | 5       | Tentative ID        | MW               | MolForm  | CAS#        | Qual |
| 1       |         | n-Hexadecanoic acid | 256              | C16H32O2 | 000057-10-3 | 96   |
| 2       |         | Tetradecanoic acid  | 228              | C14H28O2 | 000544-63-8 | 90   |
| 3       |         | Tridecanoic acid    | 214              | C13H26O2 | 000638-53-9 | 81   |
| 4       |         | Dodecanoic acid     | 200              | C12H24O2 | 000143-07-7 | 72   |
| 5       |         | Pentadecanoic acid  | 242              | C15H30O2 | 001002-84-2 | 68   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

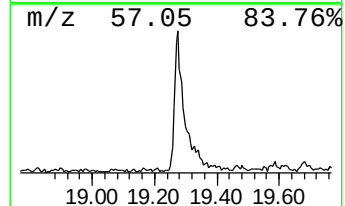
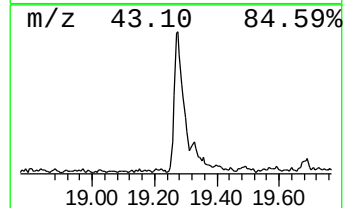
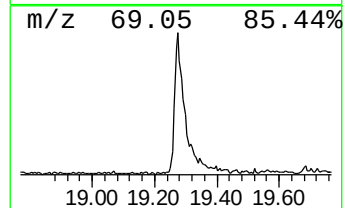
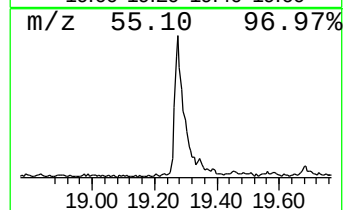
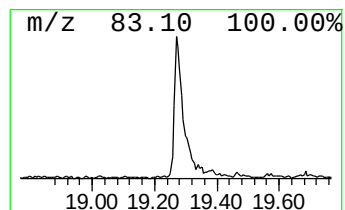
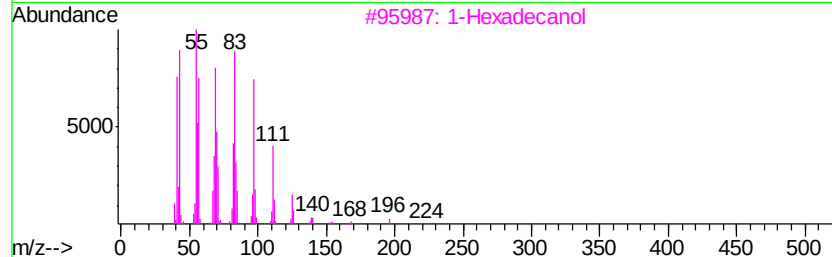
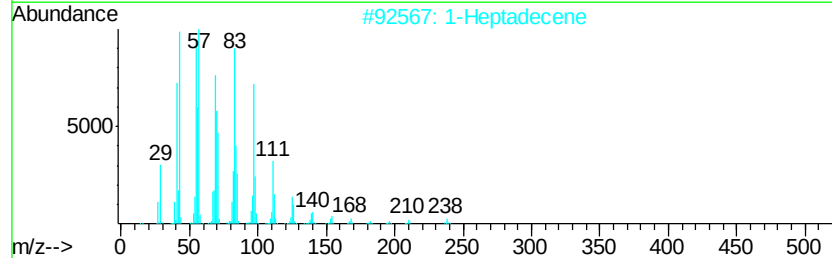
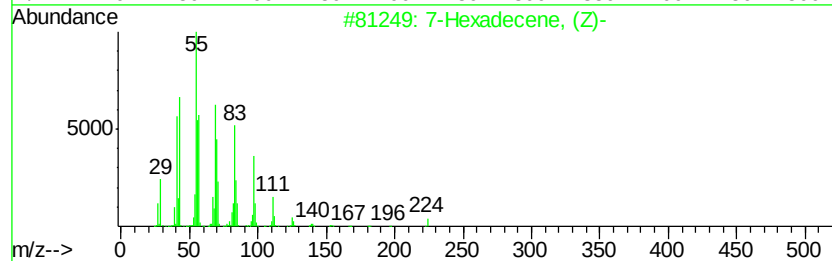
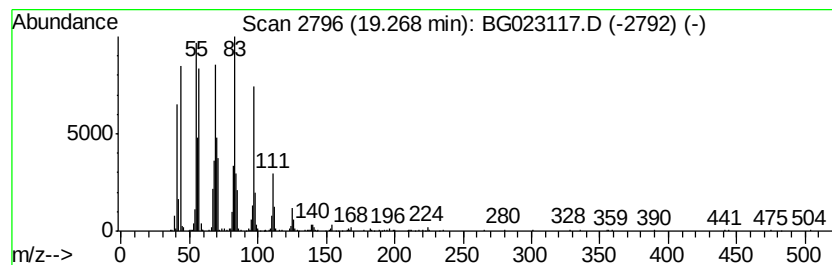
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 7-Hexadecene, (Z)- Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 19.27 | 11.36 ng | 1249790 | Phenanthrene-d10 | 17.55 |

| Hit# of | 5 | Tentative ID                       | MW  | MolForm    | CAS#        | Qual |
|---------|---|------------------------------------|-----|------------|-------------|------|
| 1       |   | 7-Hexadecene, (Z)-                 | 224 | C16H32     | 035507-09-6 | 98   |
| 2       |   | 1-Heptadecene                      | 238 | C17H34     | 006765-39-5 | 96   |
| 3       |   | 1-Hexadecanol                      | 242 | C16H34O    | 036653-82-4 | 94   |
| 4       |   | Bromoacetic acid, pentadecyl ester | 348 | C17H33BrO2 | 131143-01-6 | 94   |
| 5       |   | 9-Eicosene, (E)-                   | 280 | C20H40     | 074685-29-3 | 94   |





Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

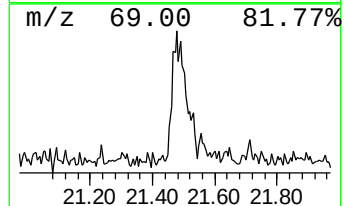
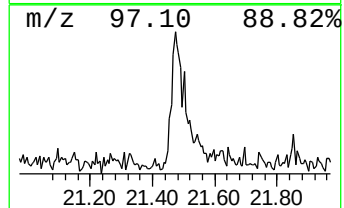
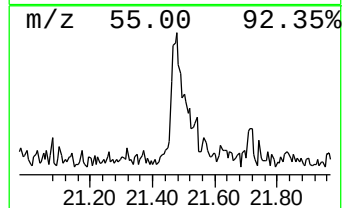
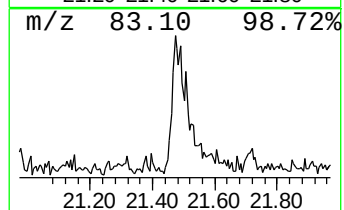
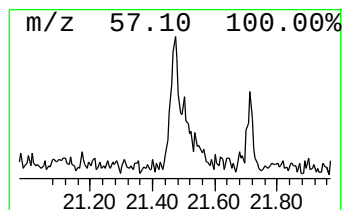
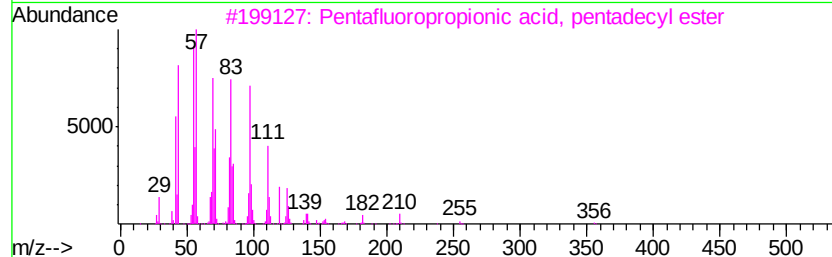
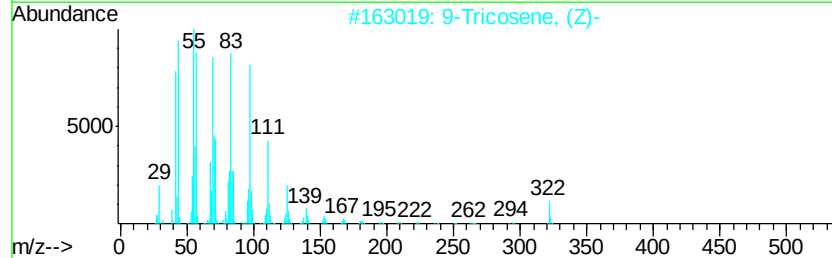
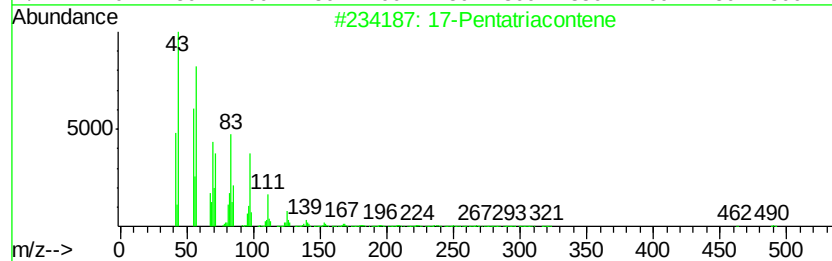
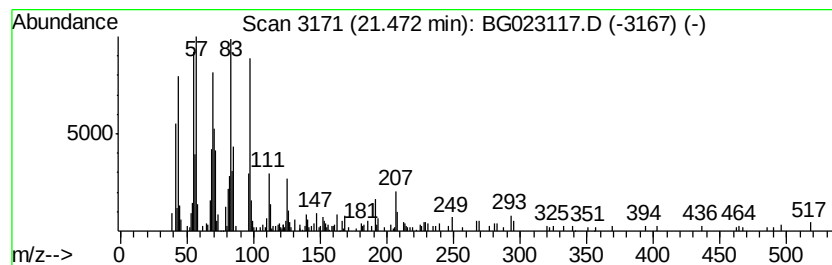
Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 17-Pentatriacontene Concentration Rank 7

| R.T.      | EstConc                             | Area   | Relative to ISTD |             | R.T.  |
|-----------|-------------------------------------|--------|------------------|-------------|-------|
| 21.47     | 2.51 ng                             | 309254 | Chrysene-d12     |             | 21.85 |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual  |
| 1         | 17-Pentatriacontene                 | 491    | C35H70           | 006971-40-0 | 91    |
| 2         | 9-Tricosene, (Z)-                   | 322    | C23H46           | 027519-02-4 | 89    |
| 3         | Pentafluoropropionic acid, penta... | 374    | C18H31F5O2       | 959092-08-1 | 74    |
| 4         | 1-Heneicosanol                      | 312    | C21H44O          | 015594-90-8 | 68    |
| 5         | 1-Hexadecanol                       | 242    | C16H34O          | 036653-82-4 | 58    |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG071516\  
Data File : BG023117.D  
Acq On : 15 Jul 2016 17:28  
Operator : UM/SJ  
Sample : H4017-17  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C5-9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG070816.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 |
| Propane, 2,2-dime... | 3.04  | 41.6    | ng    | 1502990  | 1                     | 8.15  | 722471  | 20.0 |
| unknown5.23          | 5.23  | 10.1    | ng    | 363761   | 1                     | 8.15  | 722471  | 20.0 |
| unknown7.68          | 7.68  | 85.7    | ng    | 3095090  | 1                     | 8.15  | 722471  | 20.0 |
| Cyclotetradecane     | 17.91 | 3.9     | ng    | 430238   | 4                     | 17.55 | 2200690 | 20.0 |
| n-Hexadecanoic acid  | 18.42 | 2.0     | ng    | 222145   | 4                     | 17.55 | 2200690 | 20.0 |
| 7-Hexadecene, (Z)-   | 19.27 | 11.4    | ng    | 1249790  | 4                     | 17.55 | 2200690 | 20.0 |
| 17-Pentatriacontene  | 21.47 | 2.5     | ng    | 309254   | 5                     | 21.85 | 2461580 | 20.0 |