

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
 Data File : BG016162.D  
 Acq On : 27 Mar 2015 5:04  
 Operator : TP/IZ  
 Sample : G1653-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 FD8

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-BG032515.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         | 2.793       | 14            | 18          | 32           | rVB      | 763522         | 833569        | 36.26%          | 4.808%        |
| 2         | 2.934       | 37            | 42          | 51           | rVV2     | 71816          | 128624        | 5.60%           | 0.742%        |
| 3         | 3.005       | 51            | 54          | 60           | rVB      | 63925          | 74942         | 3.26%           | 0.432%        |
| 4         | 4.920       | 373           | 380         | 389          | rVB      | 91690          | 127976        | 5.57%           | 0.738%        |
| 5         | 5.378       | 452           | 458         | 466          | rBV      | 733899         | 1054634       | 45.88%          | 6.083%        |
| 6         | 6.964       | 721           | 728         | 737          | rVB      | 753518         | 1177816       | 51.24%          | 6.793%        |
| 7         | 7.322       | 782           | 789         | 796          | rBV      | 769394         | 1219712       | 53.06%          | 7.035%        |
| 8         | 7.381       | 796           | 799         | 809          | rVB      | 24796          | 34448         | 1.50%           | 0.199%        |
| 9         | 7.569       | 825           | 831         | 838          | rVB      | 45963          | 71366         | 3.10%           | 0.412%        |
| 10        | 7.792       | 862           | 869         | 875          | rBV      | 184726         | 287806        | 12.52%          | 1.660%        |
| 11        | 8.109       | 916           | 923         | 931          | rBV      | 541958         | 860691        | 37.44%          | 4.964%        |
| 12        | 8.944       | 1057          | 1065        | 1074         | rBV      | 425863         | 700930        | 30.49%          | 4.043%        |
| 13        | 10.577      | 1336          | 1343        | 1350         | rVB      | 258498         | 436641        | 19.00%          | 2.518%        |
| 14        | 13.038      | 1755          | 1762        | 1769         | rBV      | 1095706        | 1589984       | 69.17%          | 9.170%        |
| 15        | 13.872      | 1895          | 1904        | 1912         | rBV      | 36671          | 52169         | 2.27%           | 0.301%        |
| 16        | 14.413      | 1990          | 1996        | 2004         | rVB2     | 407148         | 639531        | 27.82%          | 3.689%        |
| 17        | 15.740      | 2214          | 2222        | 2228         | rVB      | 25659          | 37828         | 1.65%           | 0.218%        |
| 18        | 15.899      | 2241          | 2249        | 2261         | rBV      | 991604         | 1379594       | 60.02%          | 7.957%        |
| 19        | 16.122      | 2283          | 2287        | 2290         | rBV      | 20770          | 26751         | 1.16%           | 0.154%        |
| 20        | 17.150      | 2456          | 2462        | 2469         | rBV2     | 542039         | 793707        | 34.53%          | 4.578%        |
| 21        | 17.591      | 2532          | 2537        | 2541         | rBV2     | 17882          | 26088         | 1.13%           | 0.150%        |
| 22        | 18.043      | 2609          | 2614        | 2621         | rBV      | 91243          | 141484        | 6.16%           | 0.816%        |
| 23        | 19.324      | 2828          | 2832        | 2841         | rBV3     | 36677          | 60472         | 2.63%           | 0.349%        |
| 24        | 19.582      | 2872          | 2876        | 2882         | rBV      | 95505          | 132059        | 5.75%           | 0.762%        |
| 25        | 19.776      | 2903          | 2909        | 2917         | rBV      | 1811379        | 2298661       | 100.00%         | 13.258%       |
| 26        | 20.387      | 3009          | 3013        | 3019         | rBV      | 16490          | 24794         | 1.08%           | 0.143%        |
| 27        | 21.033      | 3118          | 3123        | 3134         | rBV      | 210594         | 304989        | 13.27%          | 1.759%        |
| 28        | 21.321      | 3167          | 3172        | 3180         | rBV      | 732214         | 915498        | 39.83%          | 5.280%        |
| 29        | 22.508      | 3372          | 3374        | 3380         | rVB3     | 20242          | 24837         | 1.08%           | 0.143%        |
| 30        | 23.606      | 3553          | 3561        | 3569         | rBV2     | 452666         | 885200        | 38.51%          | 5.105%        |
| 31        | 26.167      | 3987          | 3997        | 4010         | rVB2     | 266103         | 792416        | 34.47%          | 4.570%        |
| 32        | 26.702      | 4081          | 4088        | 4097         | rVB2     | 38192          | 98367         | 4.28%           | 0.567%        |
| 33        | 27.483      | 4216          | 4221        | 4231         | rVB2     | 35596          | 104583        | 4.55%           | 0.603%        |

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ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Integration Parameters: rteint.p  
Integrator: RTE  
Smoothing : ON Filtering: 5  
Sampling : 1 Min Area: 3 % of largest Peak  
Start Thrs: 0.2 Max Peaks: 100  
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
Peak separation: 5

Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

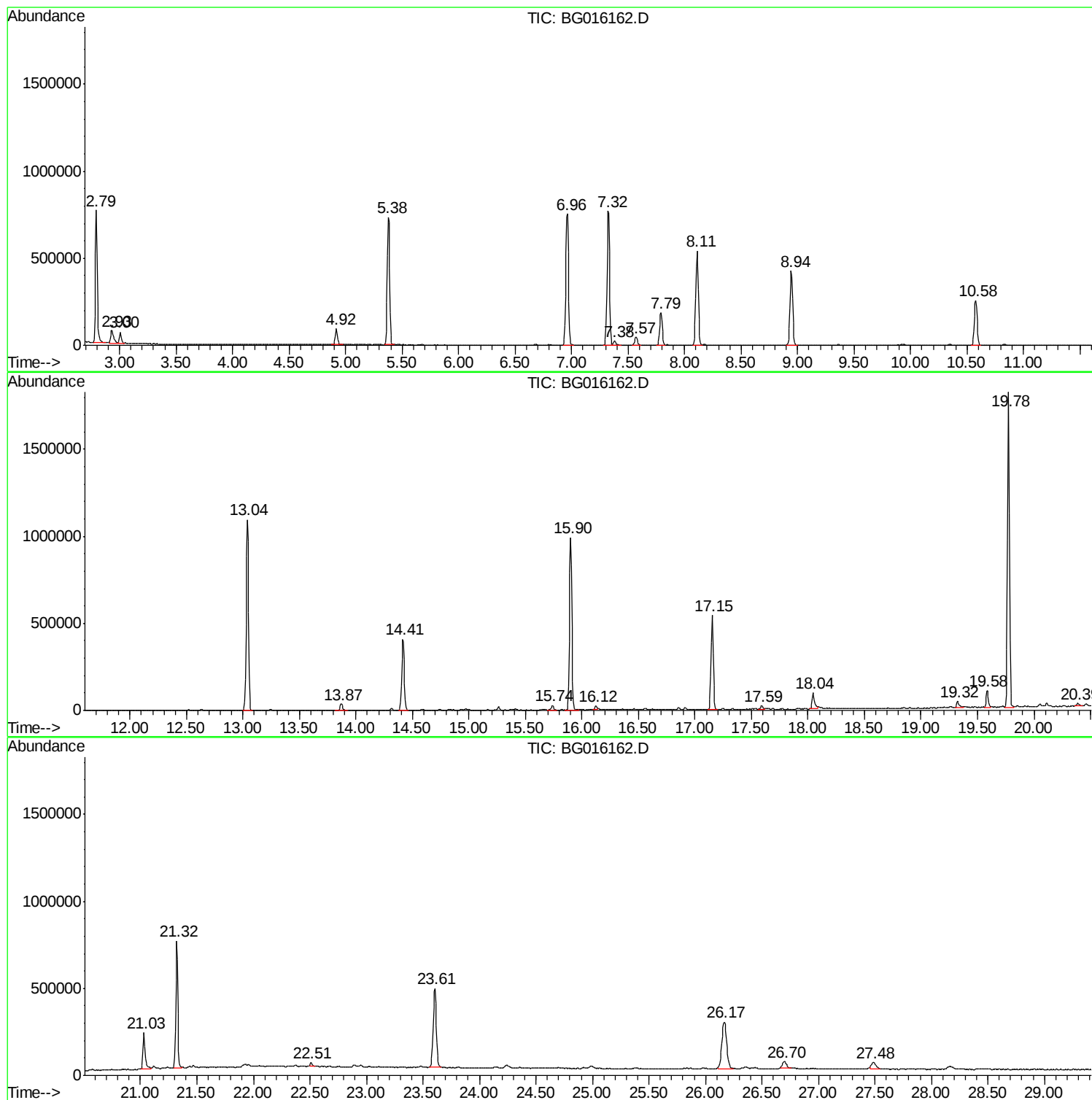
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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
FD8

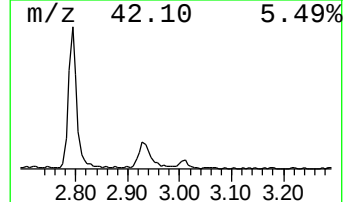
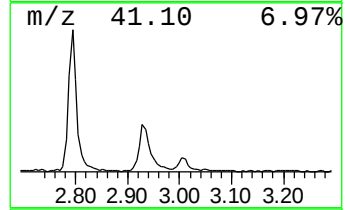
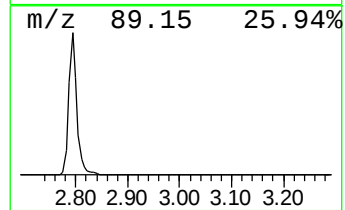
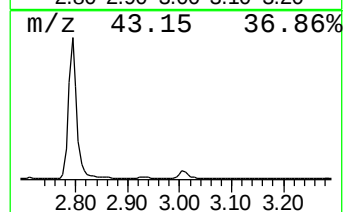
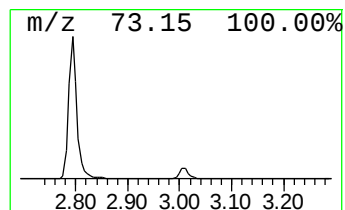
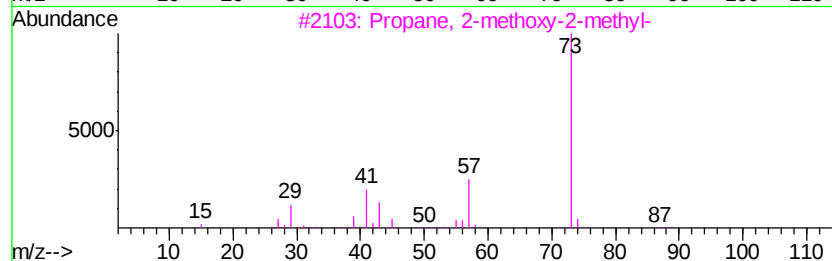
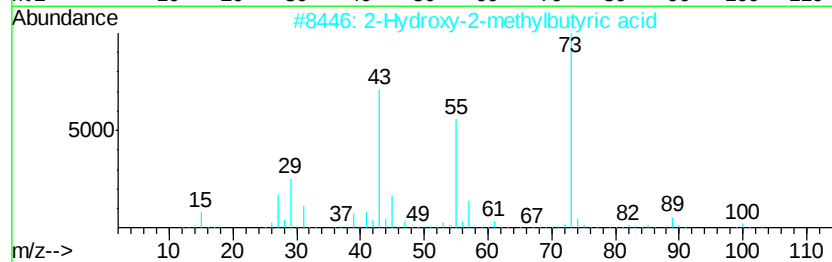
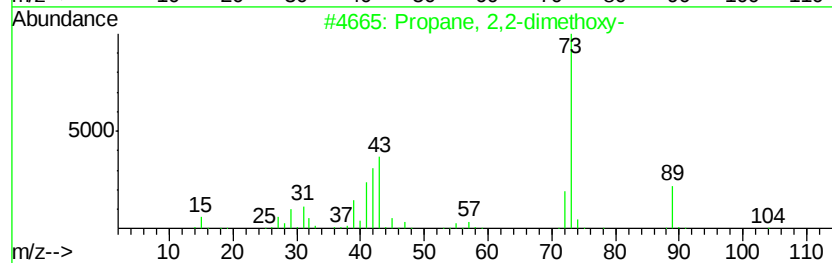
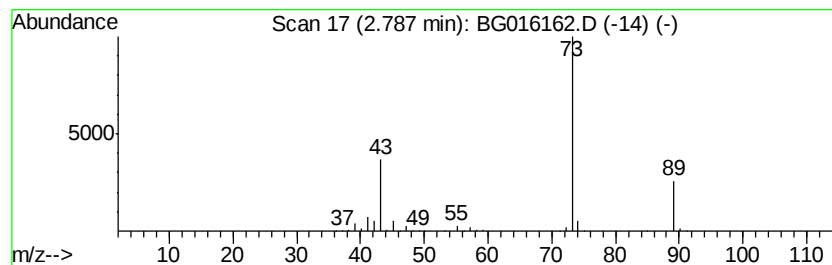
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown2.79 Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 2.79 | 57.93 ng | 833569 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Propane, 2,2-dimethoxy-        | 104 | C5H12O2 | 000077-76-9 | 45   |
| 2    |    | 2-Hydroxy-2-methylbutyric acid | 118 | C5H10O3 | 003739-30-8 | 9    |
| 3    |    | Propane, 2-methoxy-2-methyl-   | 88  | C5H12O  | 001634-04-4 | 9    |
| 4    |    | N-Ethylformamide               | 73  | C3H7NO  | 000627-45-2 | 7    |
| 5    |    | Methane, isothiocyanato-       | 73  | C2H3NS  | 000556-61-6 | 7    |



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Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

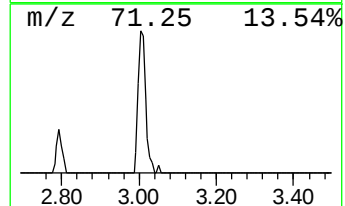
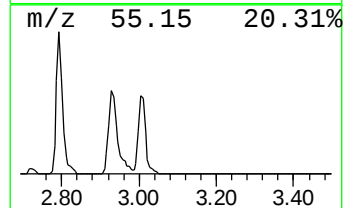
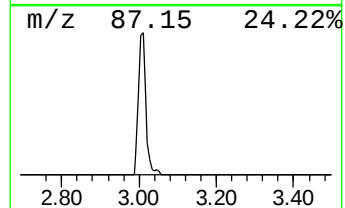
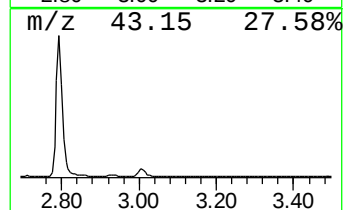
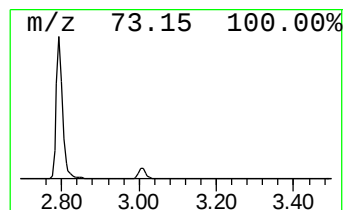
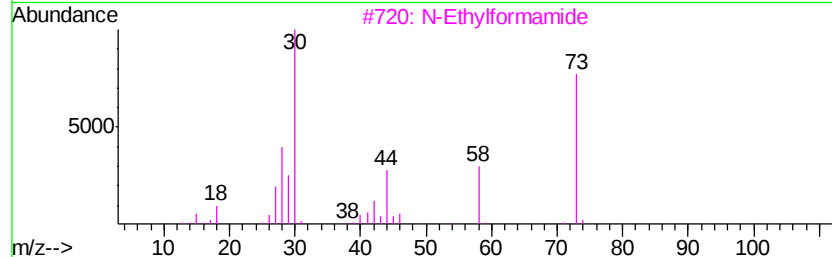
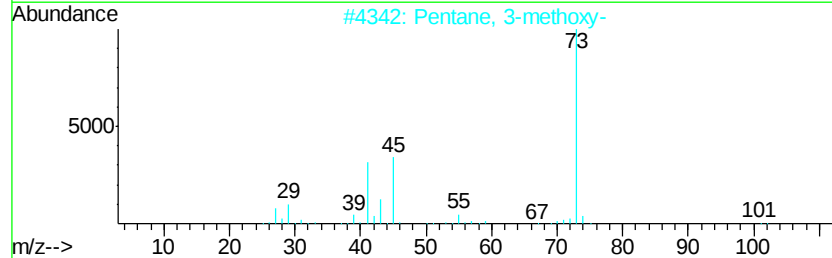
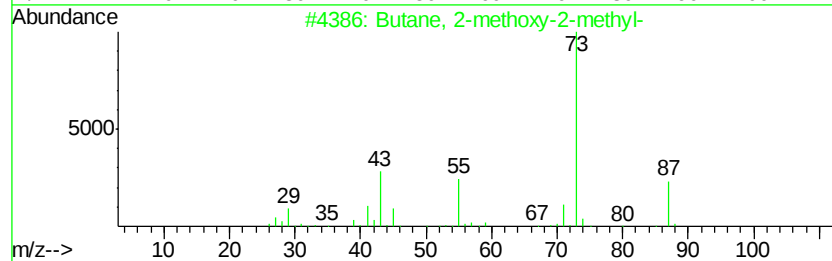
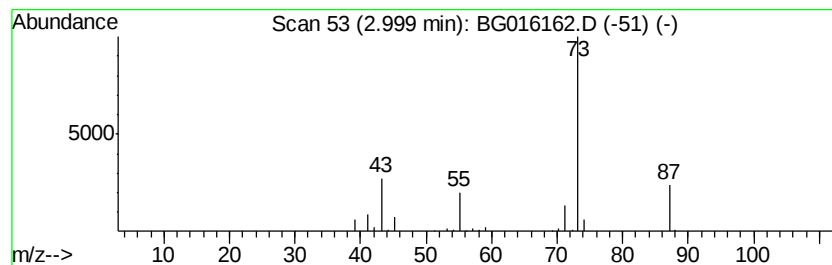
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 8

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 3.00 | 5.21 ng | 74942 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|---------|---|-----------------------------|-----|---------|-------------|------|
| 1       |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 50   |
| 2       |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 12   |
| 3       |   | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 4       |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 9    |
| 5       |   | 3-Pentanol, 2,4-dimethyl-   | 116 | C7H16O  | 000600-36-2 | 9    |



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Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

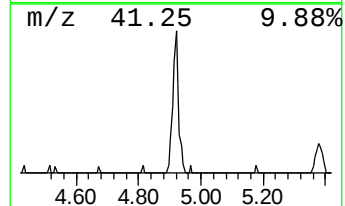
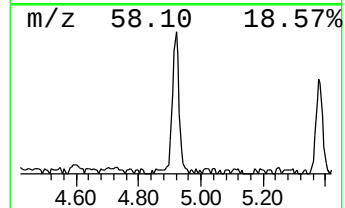
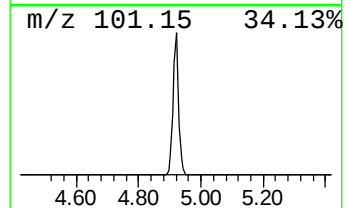
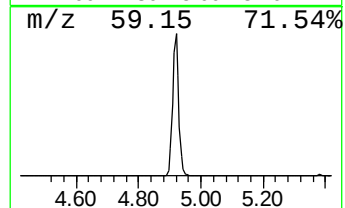
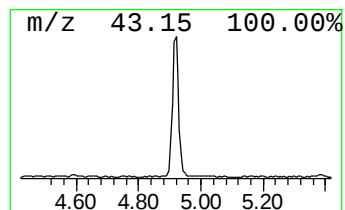
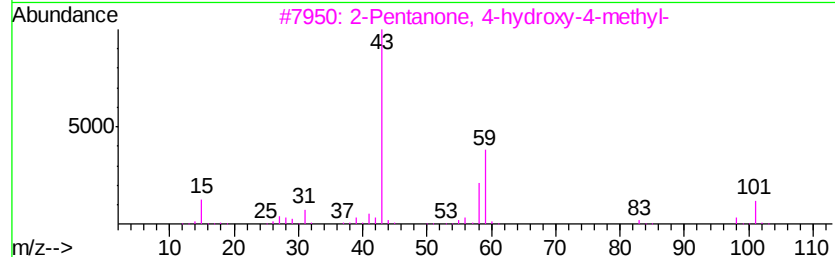
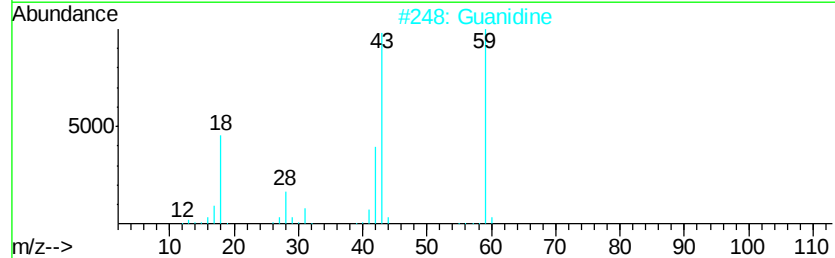
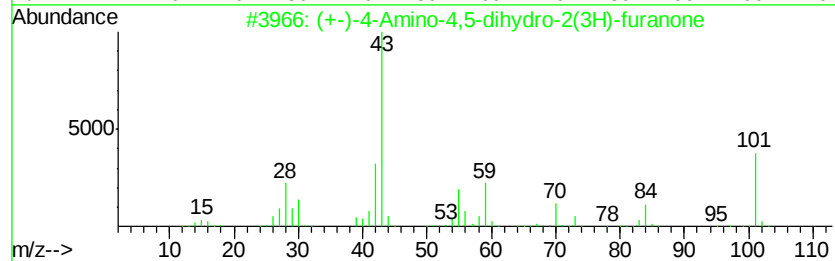
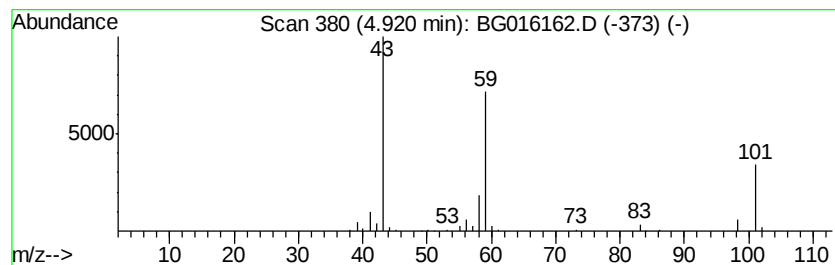
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown4.92 Concentration Rank 6

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.92 | 8.89 ng | 127976 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# | of                                  | Tentative ID | MW       | MolForm     | CAS# | Qual |
|------|-------------------------------------|--------------|----------|-------------|------|------|
| 1    | (+)-4-Amino-4,5-dihydro-2(3H)-f...  | 101          | C4H7NO2  | 016504-58-8 | 47   |      |
| 2    | Guanidine                           | 59           | CH5N3    | 000113-00-8 | 43   |      |
| 3    | 2-Pentanone, 4-hydroxy-4-methyl-    | 116          | C6H12O2  | 000123-42-2 | 40   |      |
| 4    | Acetic acid, cyano-, 1,1-dimethy... | 141          | C7H11NO2 | 001116-98-9 | 25   |      |
| 5    | 1,8-Nonanediol, 8-methyl-           | 174          | C10H22O2 | 054725-73-4 | 17   |      |



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Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

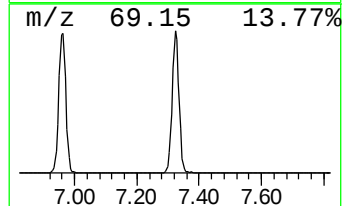
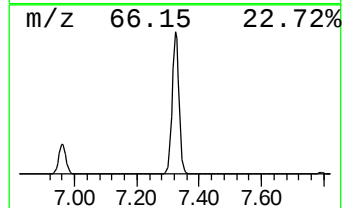
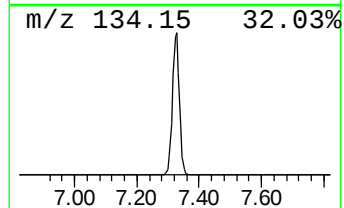
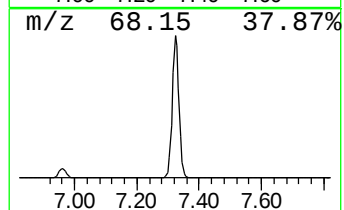
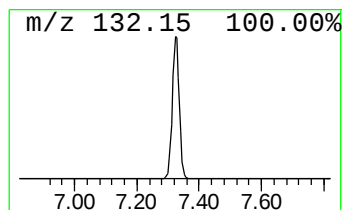
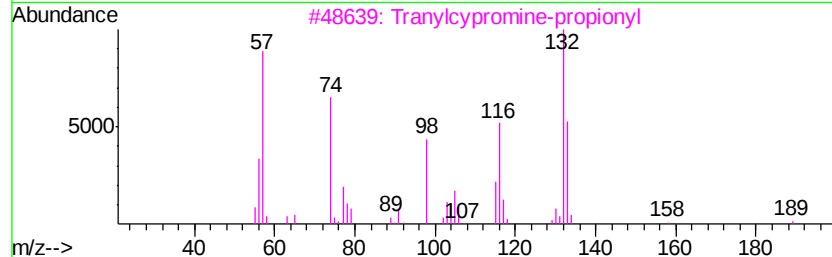
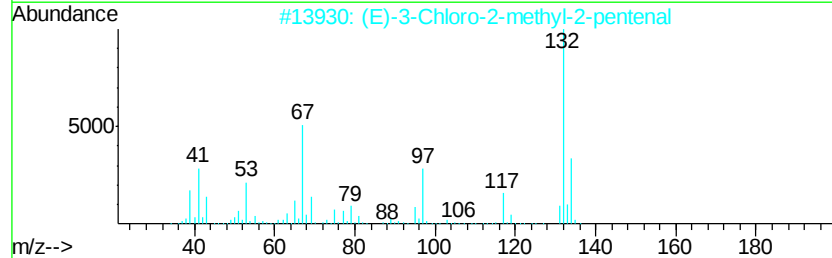
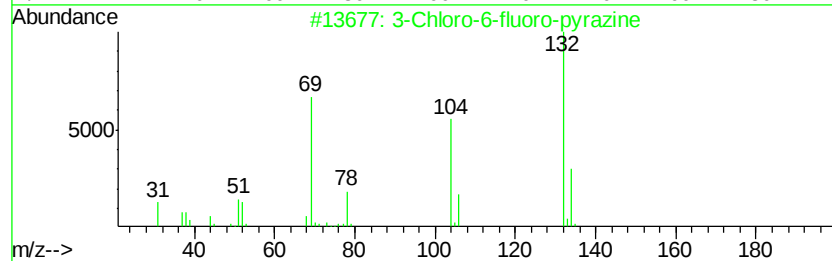
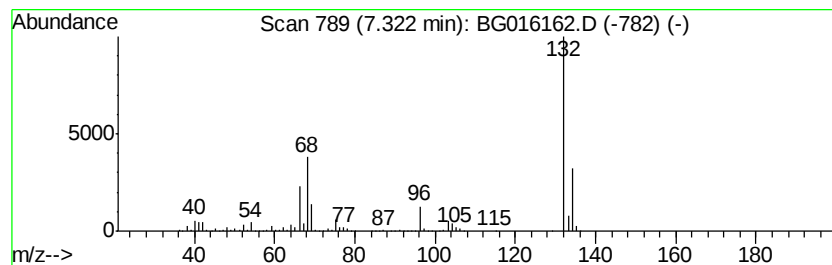
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 unknown7.32 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.32 | 84.76 ng | 1219710 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 35   |
| 3    |    | Tranylcypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2    | 022291-85-6  | 9    |
| 5    |    | 4-Chloro-2-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-00-5  | 9    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

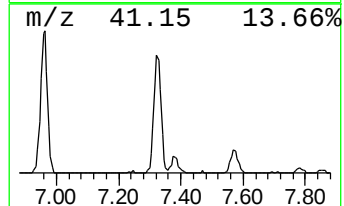
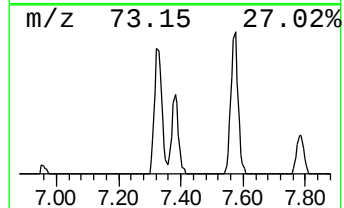
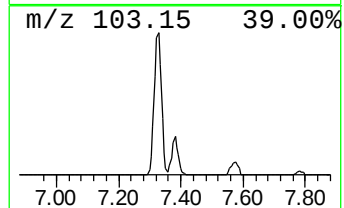
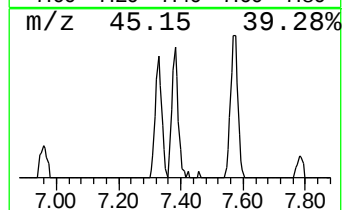
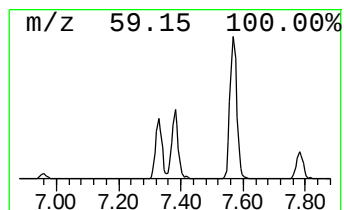
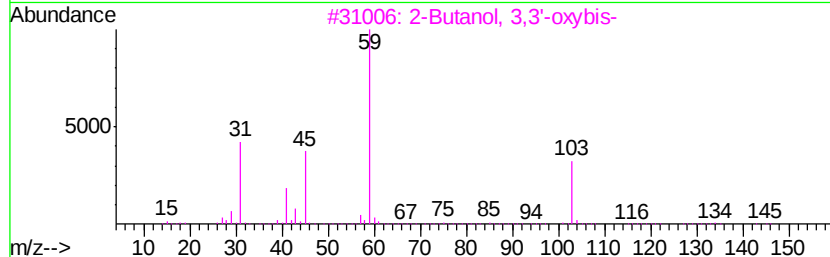
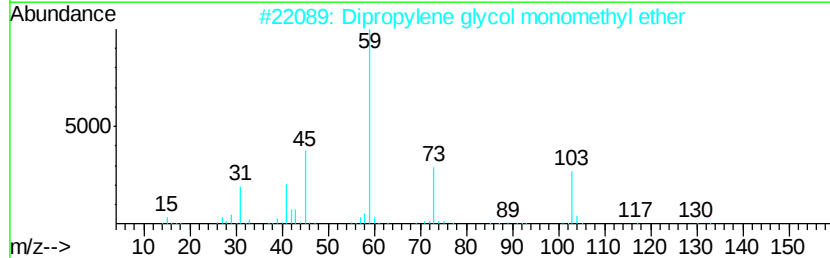
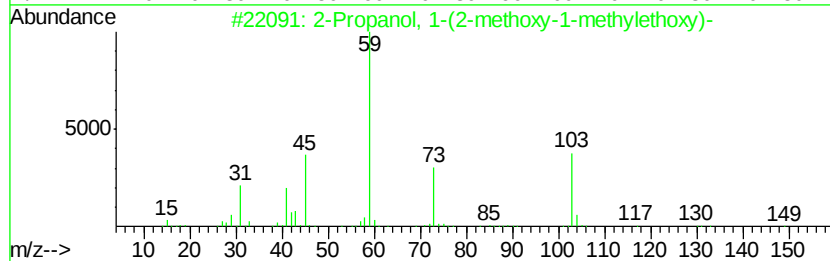
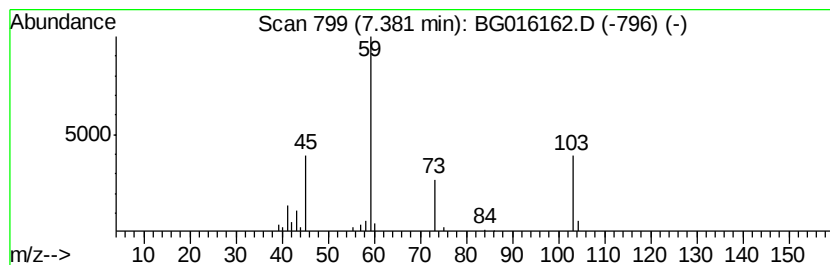
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 12

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.38 | 2.39 ng | 34448 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Propanol, 1-(2-methoxy-1-methy... | 148 | C7H16O3 | 020324-32-7 | 83   |
| 2    |    | Dipropylene glycol monomethyl ether | 148 | C7H16O3 | 034590-94-8 | 72   |
| 3    |    | 2-Butanol, 3,3'-oxybis-             | 162 | C8H18O3 | 054305-61-2 | 72   |
| 4    |    | 1-Propanol, 2-(2-hydroxypropoxy)-   | 134 | C6H14O3 | 000106-62-7 | 56   |
| 5    |    | 1-Propanol, 2,2'-oxybis-            | 134 | C6H14O3 | 000108-61-2 | 50   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016162.D  
Acq On : 27 Mar 2015 5:04  
Operator : TP/IZ  
Sample : G1653-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

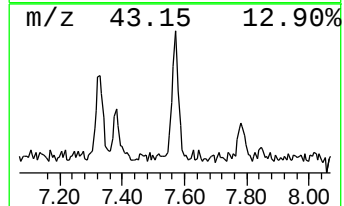
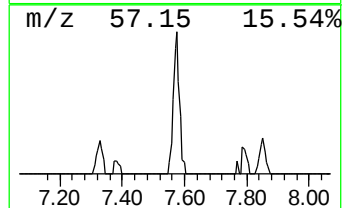
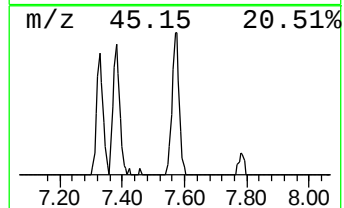
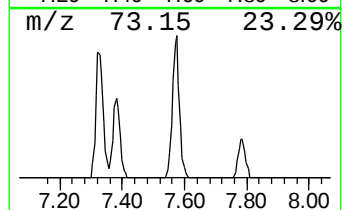
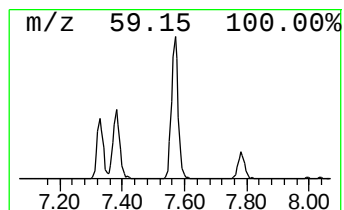
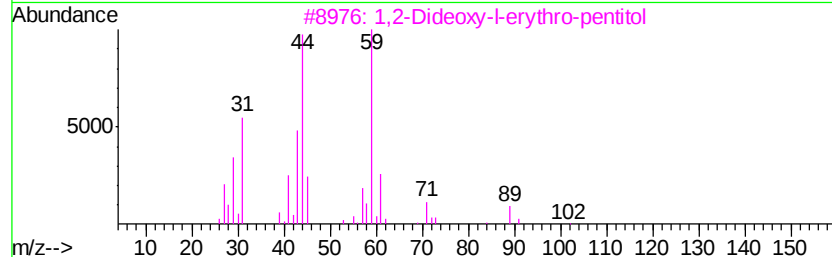
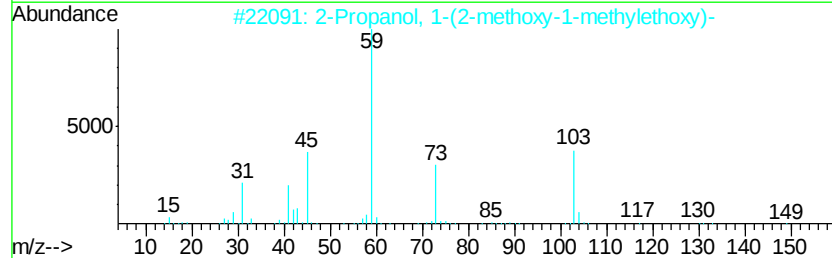
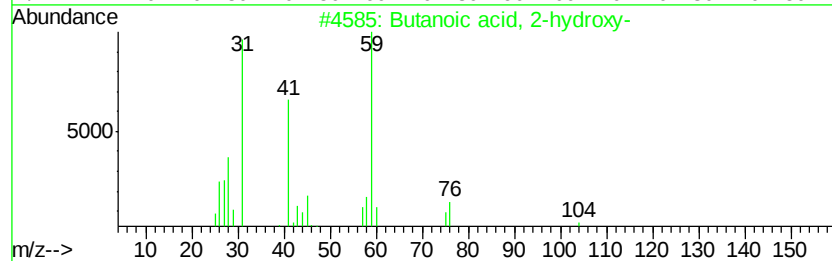
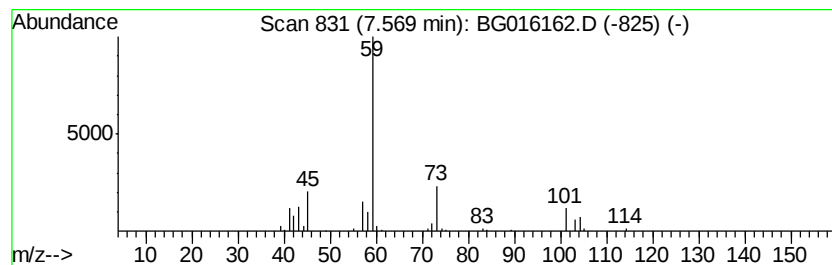
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Butanoic acid, 2-hydroxy- Concentration Rank 9

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.57 | 4.96 ng | 71366 | 1,4-Dichlorobenzene-d4 | 7.79 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|---------|---|-------------------------------------|-----|---------|--------------|------|
| 1       |   | Butanoic acid, 2-hydroxy-           | 104 | C4H8O3  | 000565-70-8  | 50   |
| 2       |   | 2-Propanol, 1-(2-methoxy-1-methy... | 148 | C7H16O3 | 020324-32-7  | 45   |
| 3       |   | 1,2-Dideoxy-1-erythro-pentitol      | 120 | C5H12O3 | 1000112-47-4 | 45   |
| 4       |   | 2-Propanol, 1-(2-methoxypropoxy)-   | 148 | C7H16O3 | 013429-07-7  | 42   |
| 5       |   | Dipropylene glycol                  | 134 | C6H14O3 | 025265-71-8  | 40   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016162.D  
Acq On : 27 Mar 2015 5:04  
Operator : TP/IZ  
Sample : G1653-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

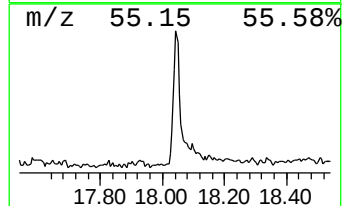
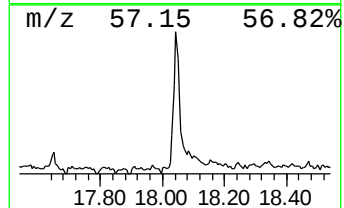
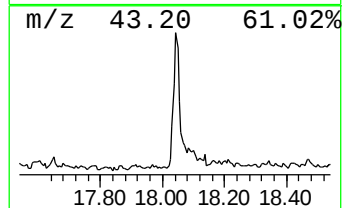
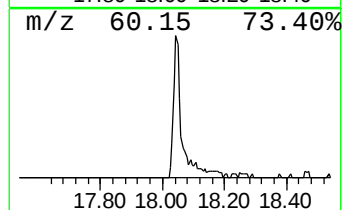
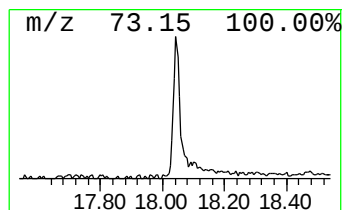
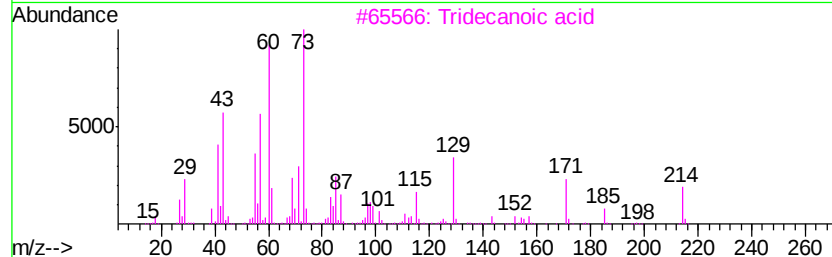
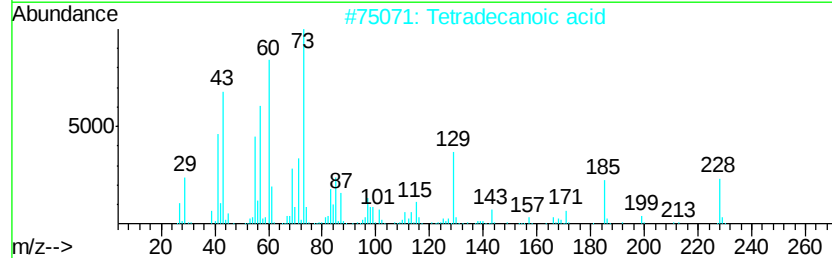
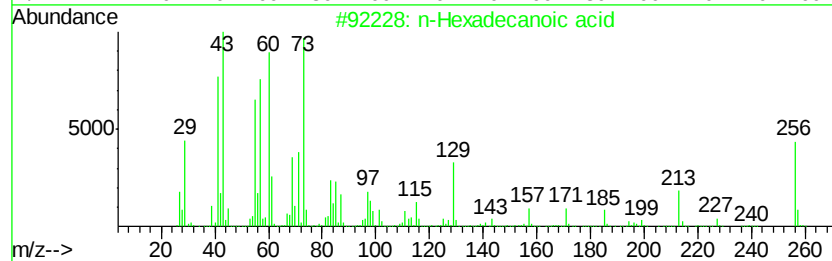
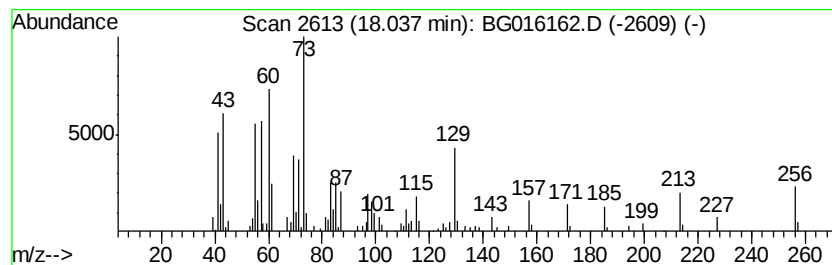
Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 n-Hexadecanoic acid Concentration Rank 10

| R.T.      | EstConc             | Area   | Relative to ISTD |             | R.T.  |
|-----------|---------------------|--------|------------------|-------------|-------|
| 18.04     | 3.57 ng             | 141484 | Phenanthrene-d10 |             | 17.15 |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual  |
| 1         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 98    |
| 2         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 96    |
| 3         | Tridecanoic acid    | 214    | C13H26O2         | 000638-53-9 | 74    |
| 4         | Dodecanoic acid     | 200    | C12H24O2         | 000143-07-7 | 47    |
| 5         | Dodecane            | 170    | C12H26           | 000112-40-3 | 11    |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016162.D  
Acq On : 27 Mar 2015 5:04  
Operator : TP/IZ  
Sample : G1653-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

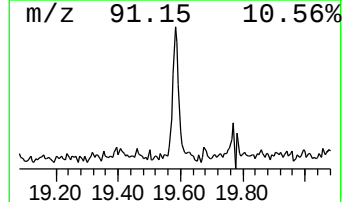
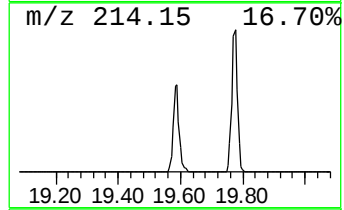
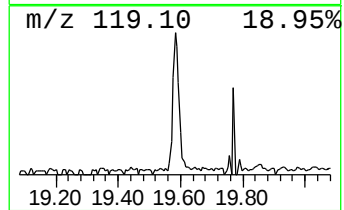
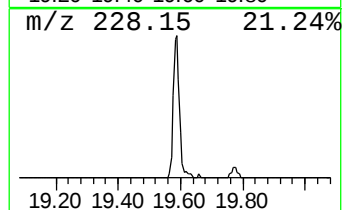
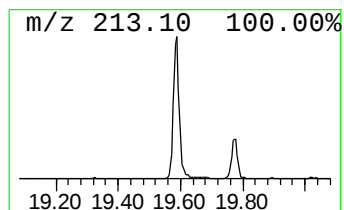
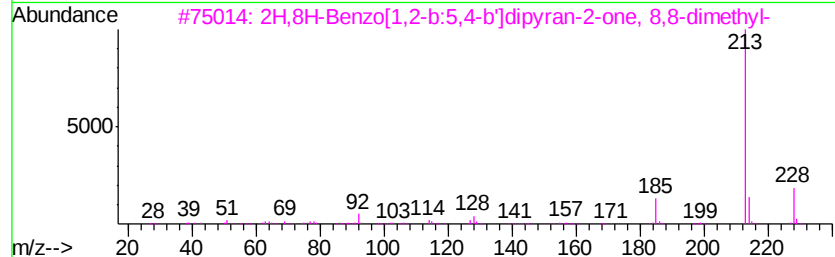
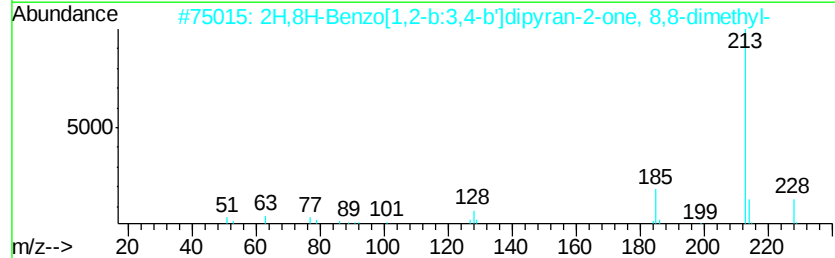
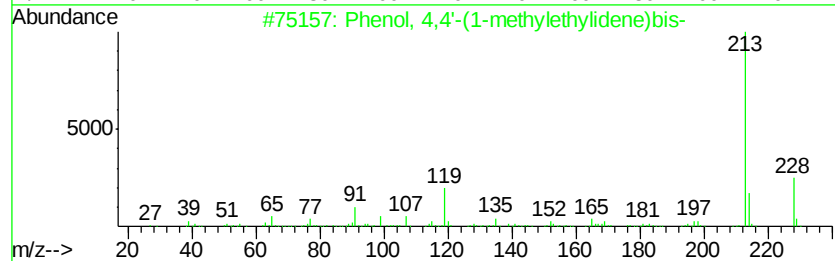
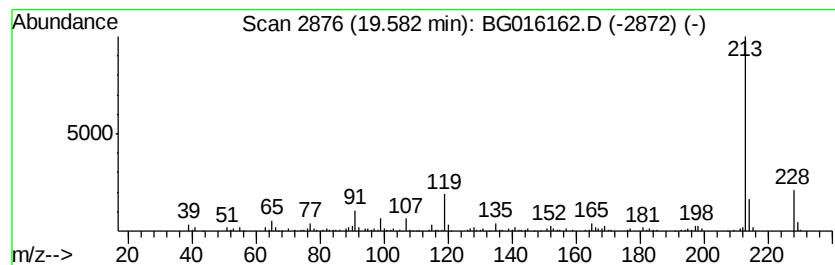
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Phenol, 4,4'-(1-methylethyl... Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.58 | 2.88 ng | 132059 | Chrysene-d12     | 21.32 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2  | 000080-05-7 | 98   |
| 2    |    | 2H,8H-Benzo[1,2-b:3,4-b']dipyran... | 228 | C14H12O3  | 000523-59-1 | 72   |
| 3    |    | 2H,8H-Benzo[1,2-b:5,4-b']dipyran... | 228 | C14H12O3  | 000553-19-5 | 64   |
| 4    |    | Carbanilic acid, p-phenyl-          | 213 | C13H11NO2 | 004474-53-7 | 47   |
| 5    |    | Phenol, 2,4'-isopropylidenedi-      | 228 | C15H16O2  | 000837-08-1 | 47   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016162.D  
Acq On : 27 Mar 2015 5:04  
Operator : TP/IZ  
Sample : G1653-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

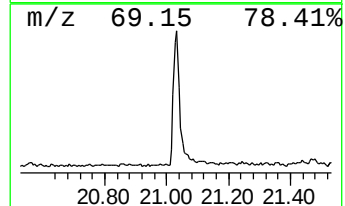
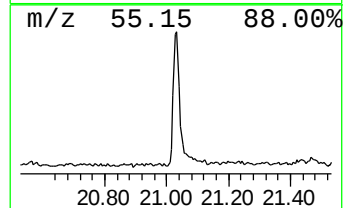
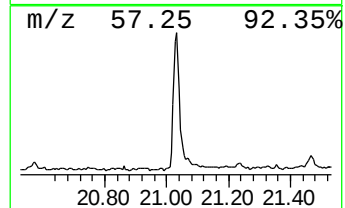
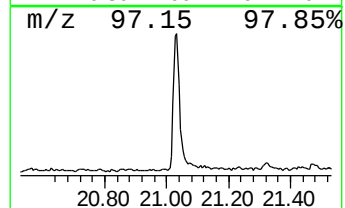
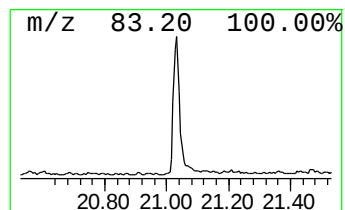
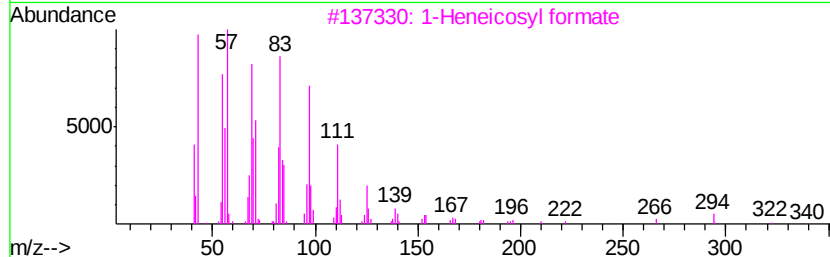
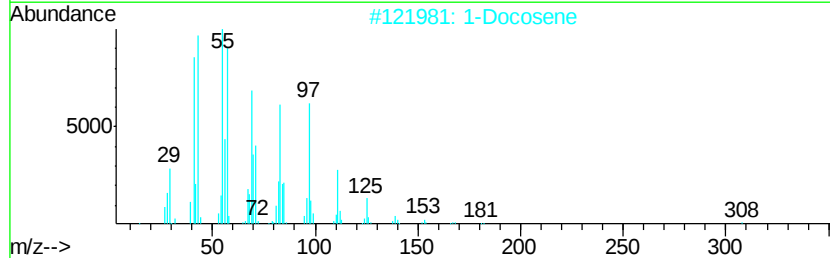
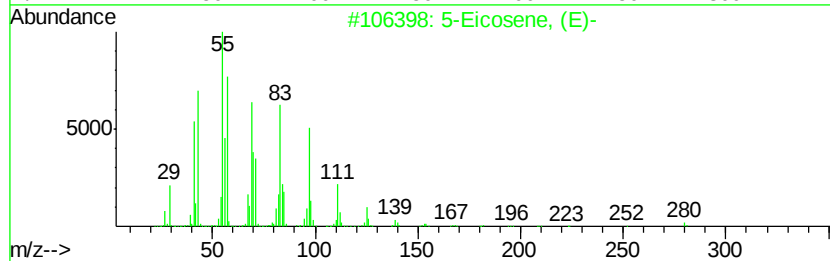
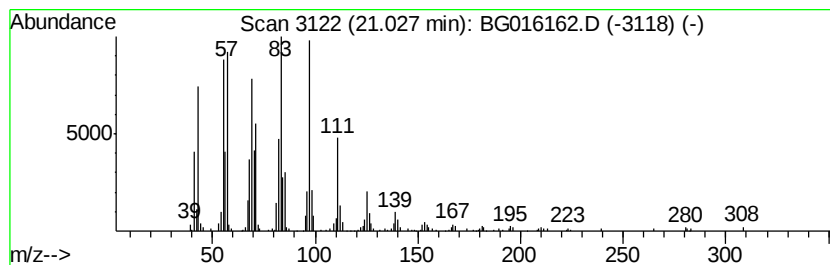
Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 5-Eicosene, (E)- Concentration Rank 7

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.            |
|---------|-------------------------------------|--------------|------------------|-----------------|
| 21.03   | 6.66 ng                             | 304989       | Chrysene-d12     | 21.32           |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual       |
| 1       | 5-Eicosene, (E)-                    |              | 280 C20H40       | 074685-30-6 99  |
| 2       | 1-Docosene                          |              | 308 C22H44       | 001599-67-3 94  |
| 3       | 1-Heneicosyl formate                |              | 340 C22H44O2     | 077899-03-7 94  |
| 4       | Cyclopentadecane                    |              | 210 C15H30       | 000295-48-7 93  |
| 5       | Dichloroacetic acid, heptadecyl ... |              | 366 C19H36Cl2O2  | 1000282-98-2 91 |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016162.D  
Acq On : 27 Mar 2015 5:04  
Operator : TP/IZ  
Sample : G1653-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

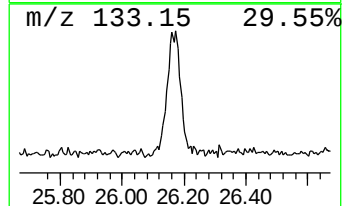
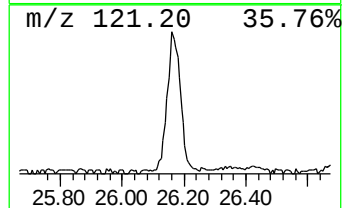
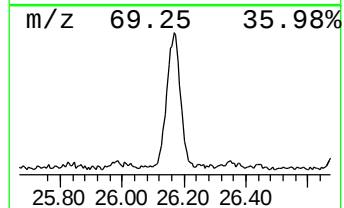
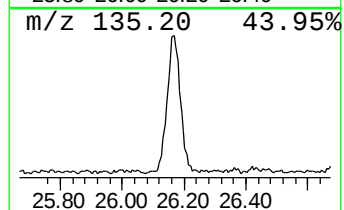
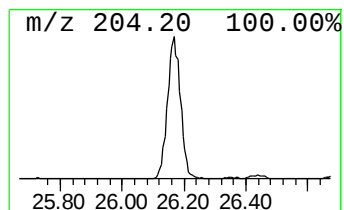
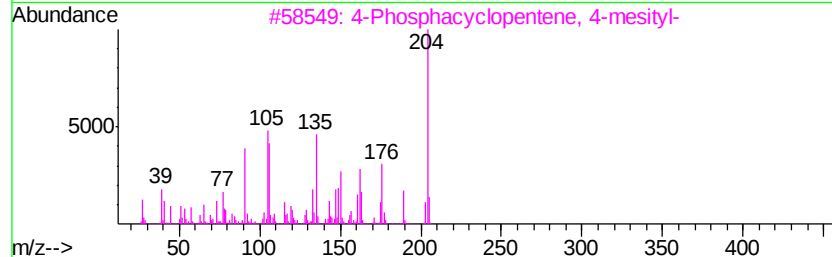
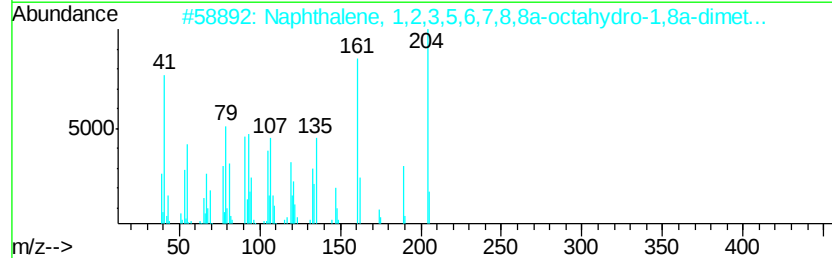
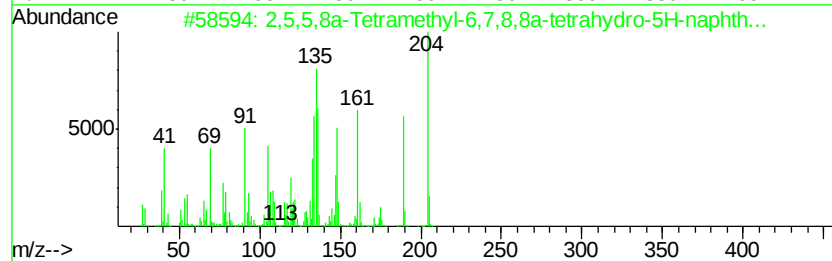
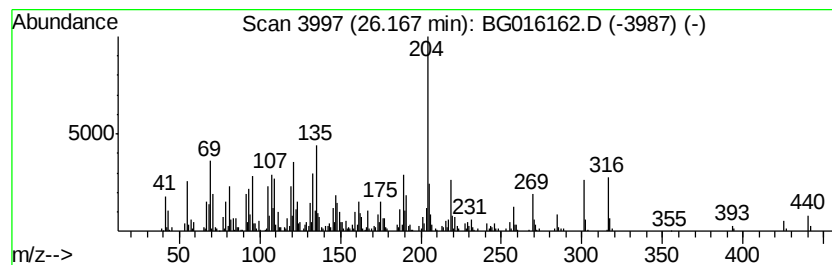
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 unknown26.17 Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 26.17 | 17.90 ng | 792416 | Perylene-d12     | 23.61 |

| Hit# | of                                  | Tentative ID | MW      | MolForm | CAS#         | Qual |
|------|-------------------------------------|--------------|---------|---------|--------------|------|
| 1    | 2,5,5,8a-Tetramethyl-6,7,8,8a-te... | 204          | C14H200 |         | 124957-09-1  | 49   |
| 2    | Naphthalene, 1,2,3,5,6,7,8,8a-oc... | 204          | C15H24  |         | 004630-07-3  | 47   |
| 3    | 4-Phosphacyclopentene, 4-mesityl-   | 204          | C13H17P |         | 214605-01-3  | 46   |
| 4    | 1,4-Dimethyl-8-isopropylidenetri... | 204          | C15H24  |         | 1000140-07-7 | 41   |
| 5    | 5H-Dibenzo[a,d]cycloheptene, 5-m... | 204          | C16H12  |         | 002975-79-3  | 27   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016162.D  
Acq On : 27 Mar 2015 5:04  
Operator : TP/IZ  
Sample : G1653-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

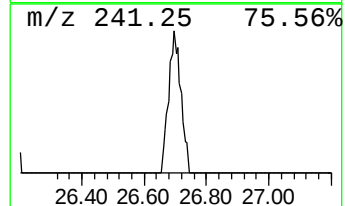
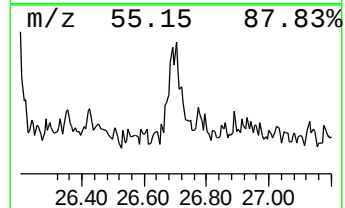
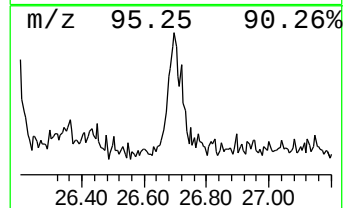
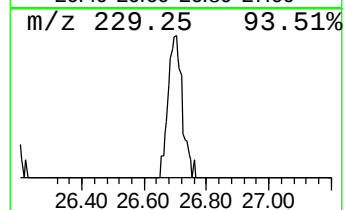
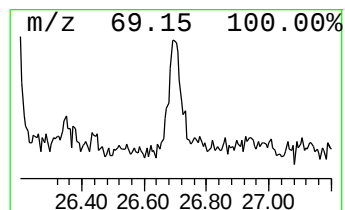
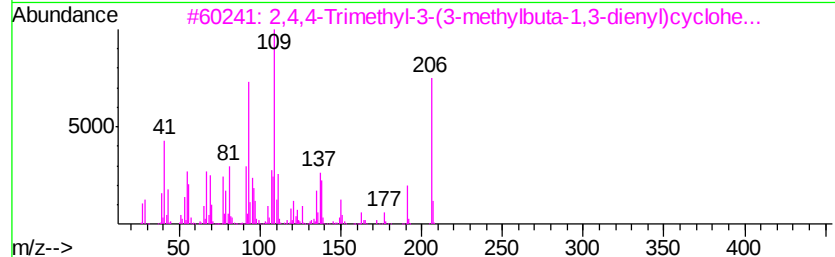
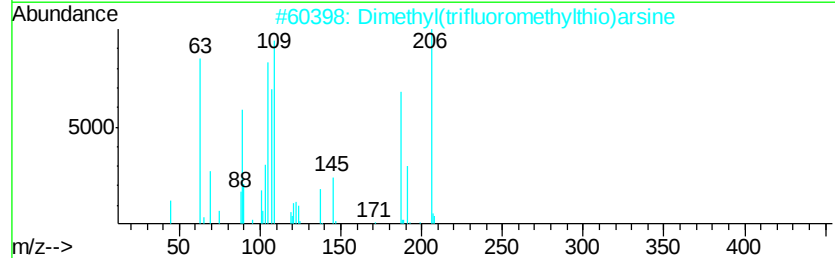
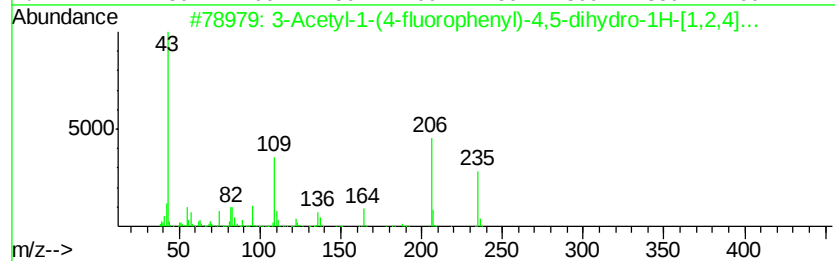
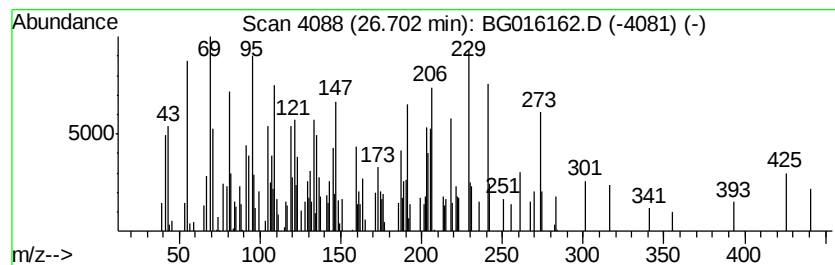
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 unknown26.70 Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 26.70 | 2.22 ng | 98367 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | 3-Acetyl-1-(4-fluorophenyl)-4,5-... | 235 | C11H10FN3O2 | 1000209-98-9 | 22   |
| 2    |    | Dimethyl(trifluoromethylthio)arsine | 206 | C3H6AsF3S   | 001959-80-4  | 14   |
| 3    |    | 2,4,4-Trimethyl-3-(3-methylbuta-... | 206 | C14H22O     | 097452-05-6  | 14   |
| 4    |    | 6,10-Dodecadien-3-ol, 3,7,11-tri... | 224 | C15H28O     | 003625-49-8  | 12   |
| 5    |    | 1-Phenanthrenecarboxylic acid, 7... | 316 | C21H32O2    | 001686-62-0  | 11   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016162.D  
Acq On : 27 Mar 2015 5:04  
Operator : TP/IZ  
Sample : G1653-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

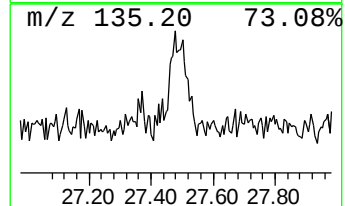
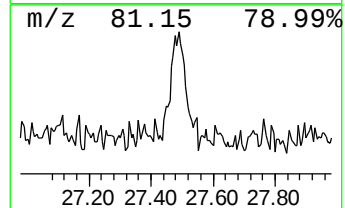
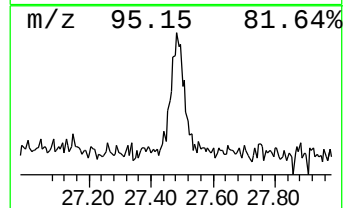
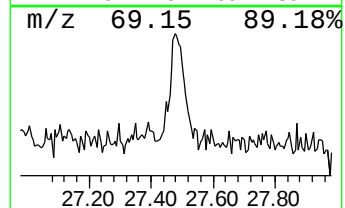
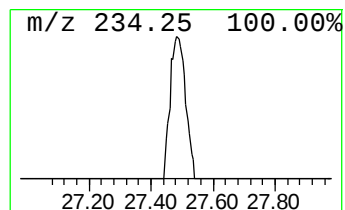
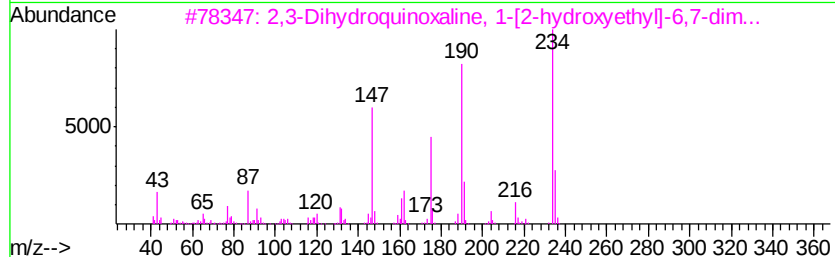
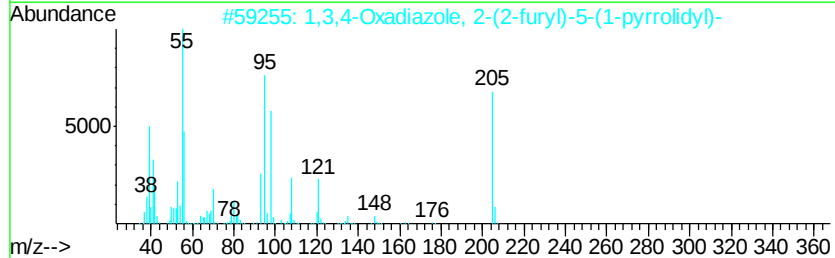
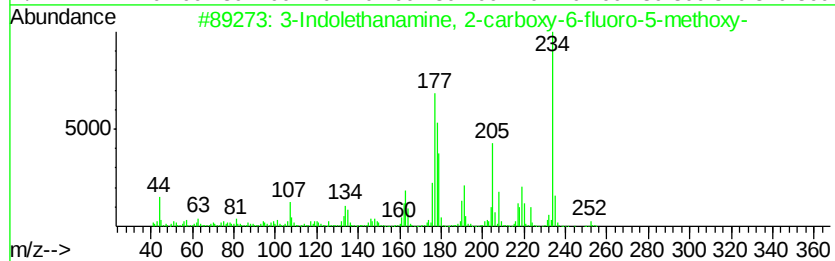
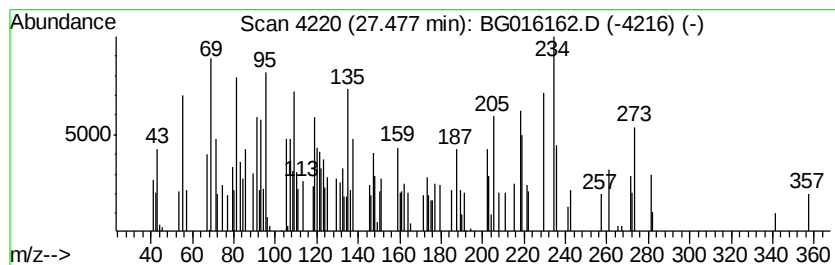
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 unknown27.48 Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 27.48 | 2.36 ng | 104583 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | 3-Indolethanamine, 2-carboxy-6-f... | 252 | C12H13FN2O3 | 062106-04-1  | 22   |
| 2    |    | 1,3,4-Oxadiazole, 2-(2-furyl)-5-... | 205 | C10H11N3O2  | 284674-14-2  | 14   |
| 3    |    | 2,3-Dihydroquinoxaline, 1-[2-hyd... | 234 | C12H14N2O3  | 1000127-85-4 | 11   |
| 4    |    | Benzoic acid, 2-hydroximinoacety... | 208 | C9H8N2O4    | 006579-46-0  | 11   |
| 5    |    | 5-Acetyl-2-methylpyridine thiose... | 208 | C9H12N4S    | 219865-93-7  | 11   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG032615\  
Data File : BG016162.D  
Acq On : 27 Mar 2015 5:04  
Operator : TP/IZ  
Sample : G1653-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
FD8

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG032515.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| unknown2.79          | 2.79  | 57.9    | ng    | 833569   | 1                     | 7.79  | 287806 | 20.0 |
| Butane, 2-methoxy... | 3.00  | 5.2     | ng    | 74942    | 1                     | 7.79  | 287806 | 20.0 |
| unknown4.92          | 4.92  | 8.9     | ng    | 127976   | 1                     | 7.79  | 287806 | 20.0 |
| unknown7.32          | 7.32  | 84.8    | ng    | 1219710  | 1                     | 7.79  | 287806 | 20.0 |
| 2-Propanol, 1-(2-... | 7.38  | 2.4     | ng    | 34448    | 1                     | 7.79  | 287806 | 20.0 |
| Butanoic acid, 2-... | 7.57  | 5.0     | ng    | 71366    | 1                     | 7.79  | 287806 | 20.0 |
| n-Hexadecanoic acid  | 18.04 | 3.6     | ng    | 141484   | 4                     | 17.15 | 793707 | 20.0 |
| Phenol, 4,4'-(1-m... | 19.58 | 2.9     | ng    | 132059   | 5                     | 21.32 | 915498 | 20.0 |
| 5-Eicosene, (E)-     | 21.03 | 6.7     | ng    | 304989   | 5                     | 21.32 | 915498 | 20.0 |
| unknown26.17         | 26.17 | 17.9    | ng    | 792416   | 6                     | 23.61 | 885200 | 20.0 |
| unknown26.70         | 26.70 | 2.2     | ng    | 98367    | 6                     | 23.61 | 885200 | 20.0 |
| unknown27.48         | 27.48 | 2.4     | ng    | 104583   | 6                     | 23.61 | 885200 | 20.0 |