

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121215\  
 Data File : BG020118.D  
 Acq On : 11 Dec 2015 22:43  
 Operator : UM/SJ  
 Sample : G4729-15  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 W-42

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-BG120215.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         | 3.120       | 43            | 48          | 57           | rBV2     | 35315          | 80639         | 3.78%           | 0.961%        |
| 2         | 3.202       | 57            | 62          | 71           | rVB      | 31765          | 53346         | 2.50%           | 0.636%        |
| 3         | 5.176       | 392           | 398         | 406          | rBV      | 32524          | 50003         | 2.34%           | 0.596%        |
| 4         | 5.646       | 471           | 478         | 485          | rBV      | 142262         | 222433        | 10.42%          | 2.652%        |
| 5         | 7.250       | 744           | 751         | 758          | rBV      | 98269          | 169991        | 7.96%           | 2.027%        |
| 6         | 7.626       | 808           | 815         | 823          | rBV      | 255699         | 436044        | 20.43%          | 5.199%        |
| 7         | 8.102       | 887           | 896         | 902          | rBV      | 68478          | 114148        | 5.35%           | 1.361%        |
| 8         | 8.425       | 943           | 951         | 960          | rBV      | 232302         | 396920        | 18.59%          | 4.733%        |
| 9         | 9.266       | 1086          | 1094        | 1102         | rBV      | 207968         | 363660        | 17.03%          | 4.336%        |
| 10        | 10.917      | 1368          | 1375        | 1382         | rBV      | 97494          | 168688        | 7.90%           | 2.011%        |
| 11        | 13.355      | 1783          | 1790        | 1797         | rVB      | 690850         | 1060112       | 49.66%          | 12.640%       |
| 12        | 14.624      | 1998          | 2006        | 2013         | rBV2     | 18977          | 38224         | 1.79%           | 0.456%        |
| 13        | 14.730      | 2017          | 2024        | 2031         | rVB2     | 175162         | 278030        | 13.02%          | 3.315%        |
| 14        | 16.211      | 2265          | 2276        | 2287         | rBV      | 659492         | 990170        | 46.38%          | 11.806%       |
| 15        | 16.416      | 2307          | 2311        | 2323         | rVB      | 13254          | 23132         | 1.08%           | 0.276%        |
| 16        | 17.468      | 2482          | 2490        | 2496         | rBV      | 255085         | 379020        | 17.75%          | 4.519%        |
| 17        | 17.814      | 2544          | 2549        | 2564         | rBV      | 49929          | 73576         | 3.45%           | 0.877%        |
| 18        | 18.337      | 2631          | 2638        | 2645         | rBV2     | 49651          | 72946         | 3.42%           | 0.870%        |
| 19        | 19.607      | 2849          | 2854        | 2863         | rBV      | 44372          | 57209         | 2.68%           | 0.682%        |
| 20        | 19.712      | 2866          | 2872        | 2877         | rVV      | 122549         | 172811        | 8.09%           | 2.060%        |
| 21        | 20.071      | 2927          | 2933        | 2939         | rBV      | 1701735        | 2134791       | 100.00%         | 25.454%       |
| 22        | 20.840      | 3060          | 3064        | 3073         | rBV3     | 18386          | 29388         | 1.38%           | 0.350%        |
| 23        | 21.211      | 3123          | 3127        | 3135         | rBV      | 62769          | 92806         | 4.35%           | 1.107%        |
| 24        | 21.739      | 3210          | 3217        | 3223         | rBV      | 296784         | 467861        | 21.92%          | 5.578%        |
| 25        | 22.909      | 3410          | 3416        | 3426         | rBV3     | 13383          | 27810         | 1.30%           | 0.332%        |
| 26        | 25.006      | 3761          | 3773        | 3784         | rBV2     | 150932         | 433216        | 20.29%          | 5.165%        |

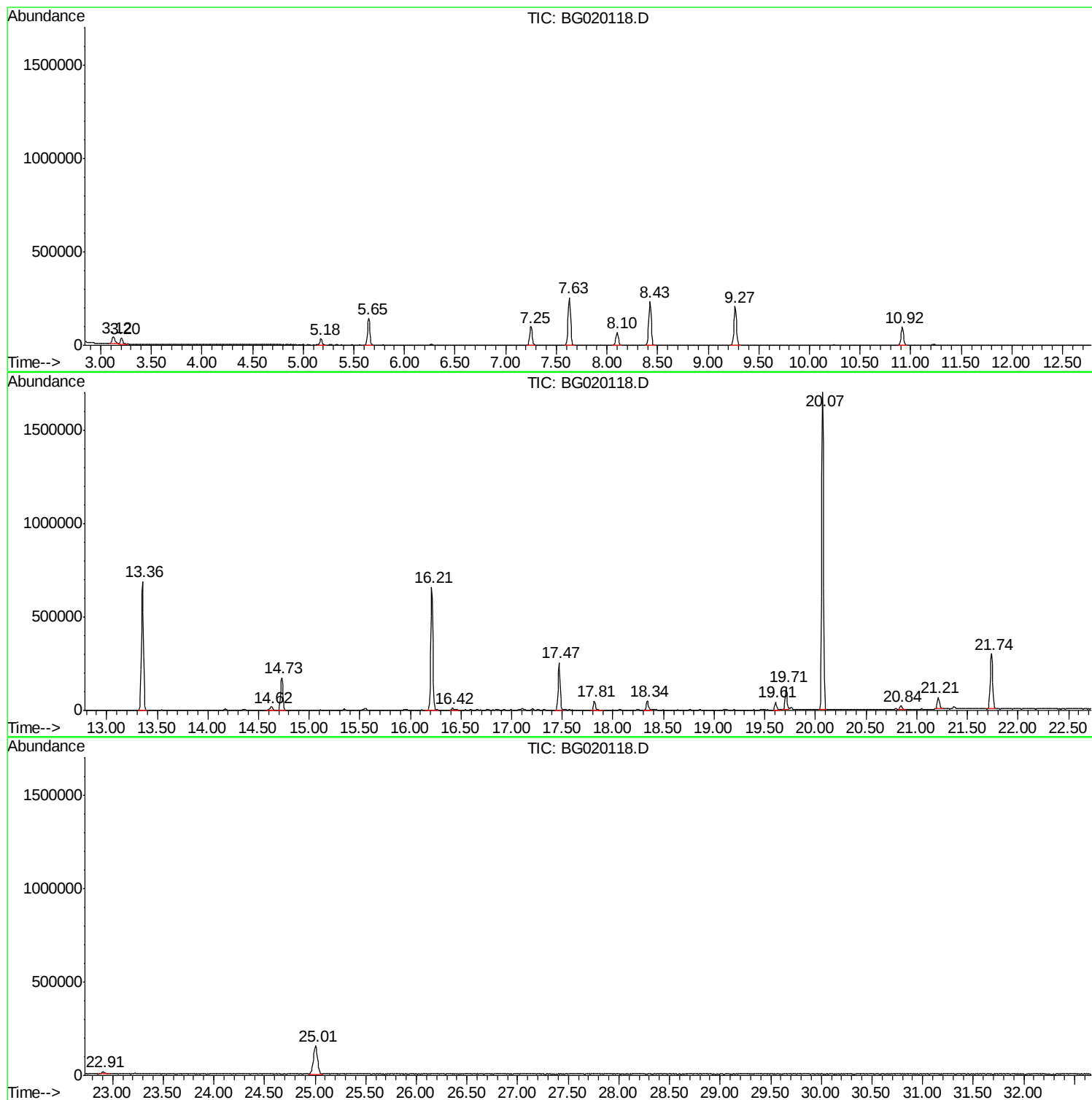
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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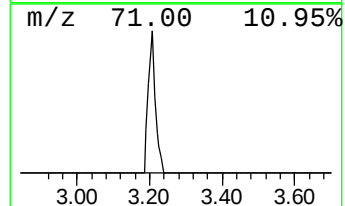
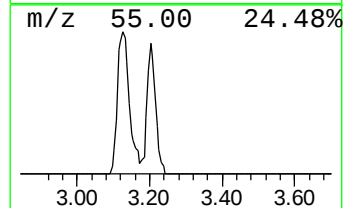
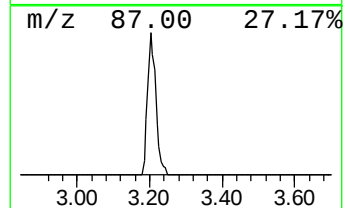
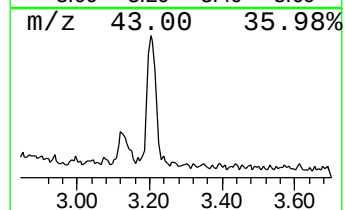
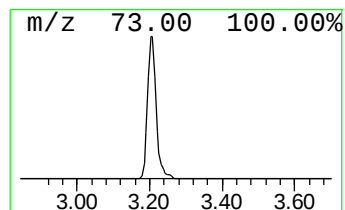
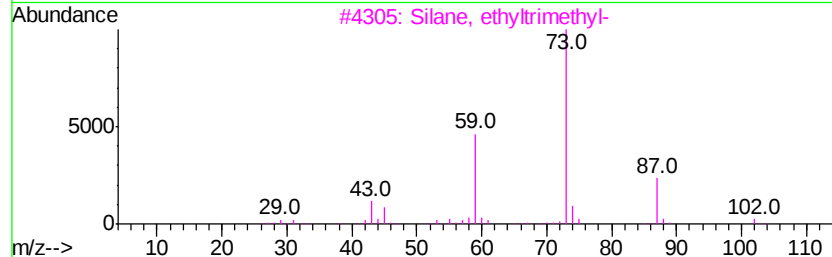
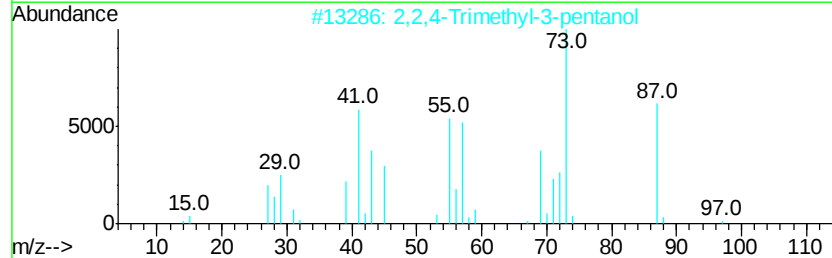
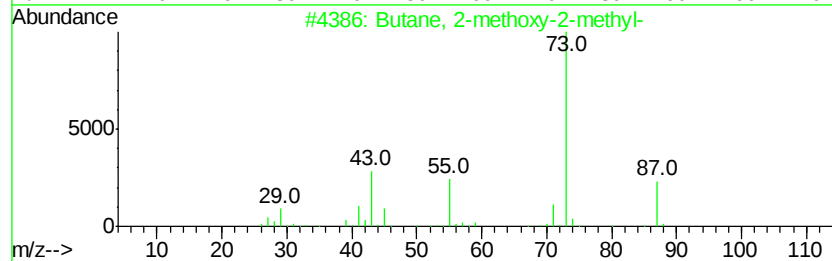
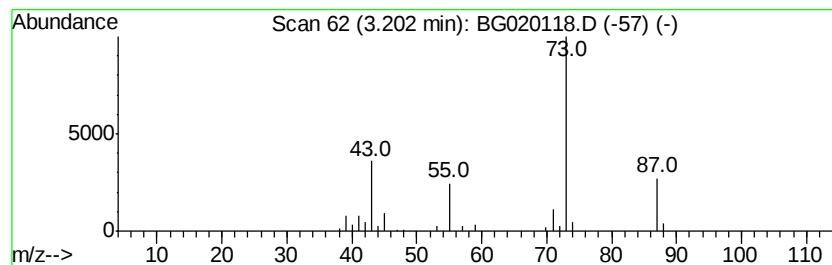
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 3.20 | 9.35 ng | 53346 | 1,4-Dichlorobenzene-d4 | 8.10 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    | Silane, ethyltrimethyl-     | 102 | C5H14Si | 003439-38-1 | 36   |
| 4    |    | 2-Methyl-5-hexen-3-ol       | 114 | C7H14O  | 032815-70-6 | 25   |
| 5    |    | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 9    |



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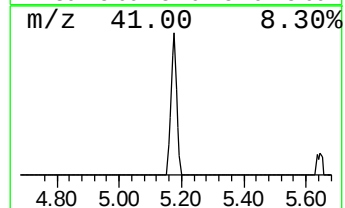
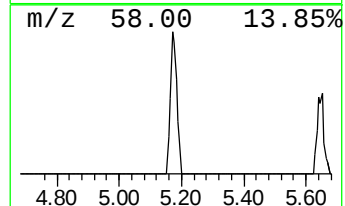
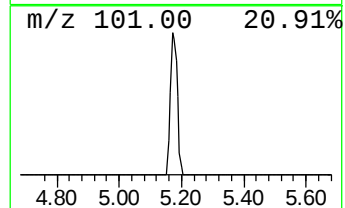
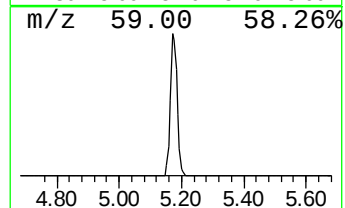
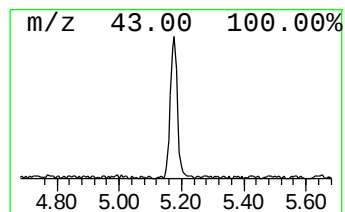
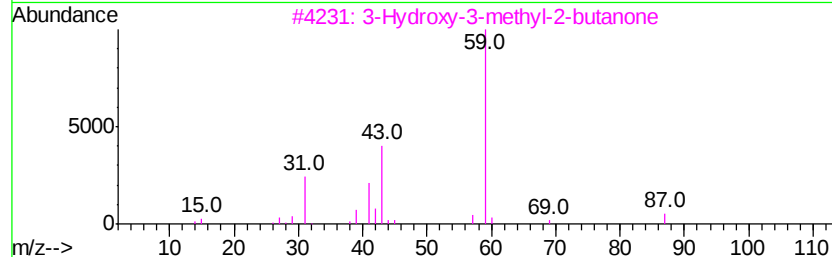
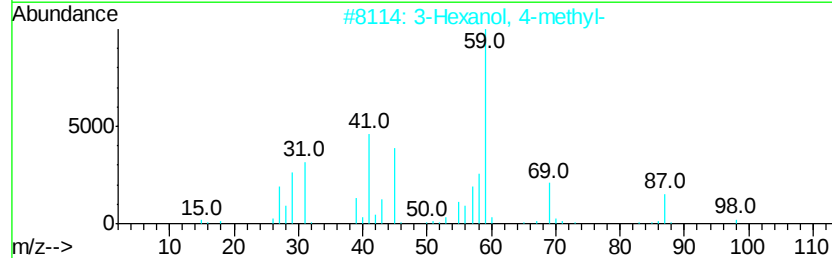
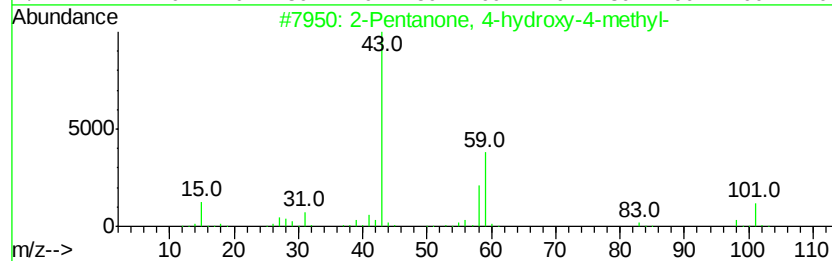
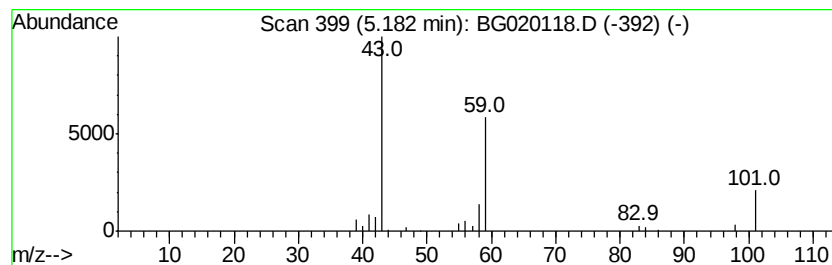
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown5.18 Concentration Rank 5

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 5.18 | 8.76 ng | 50003 | 1,4-Dichlorobenzene-d4 | 8.10 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2      | 000123-42-2 | 33      |      |      |
| 2    | 3-Hexanol, 4-methyl-             | 116 | C7H16O       | 000615-29-2 | 28      |      |      |
| 3    | 3-Hydroxy-3-methyl-2-butanone    | 102 | C5H10O2      | 000115-22-0 | 17      |      |      |
| 4    | 5-Hexen-2-one                    | 98  | C6H10O       | 000109-49-9 | 9       |      |      |
| 5    | 2,3-Butanedione, monooxime       | 101 | C4H7NO2      | 000057-71-6 | 9       |      |      |



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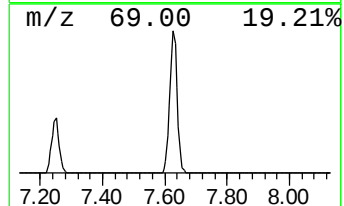
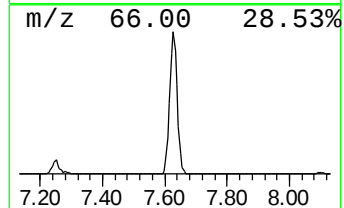
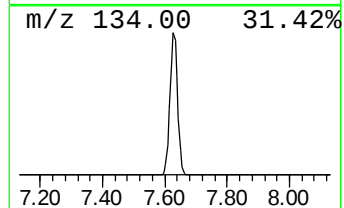
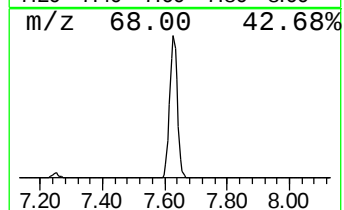
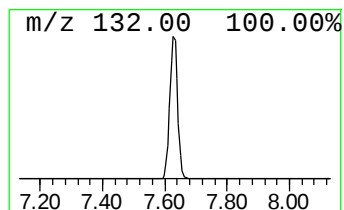
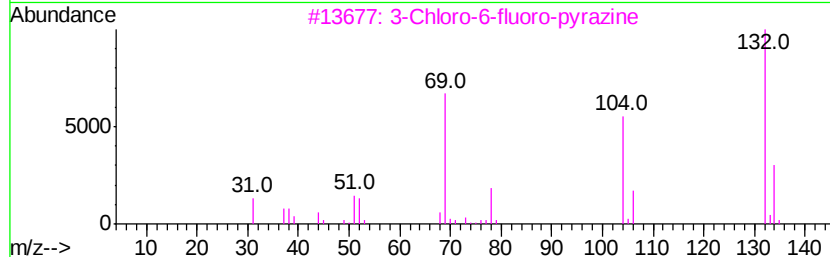
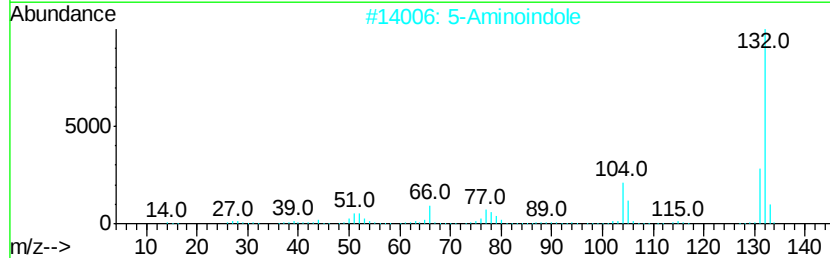
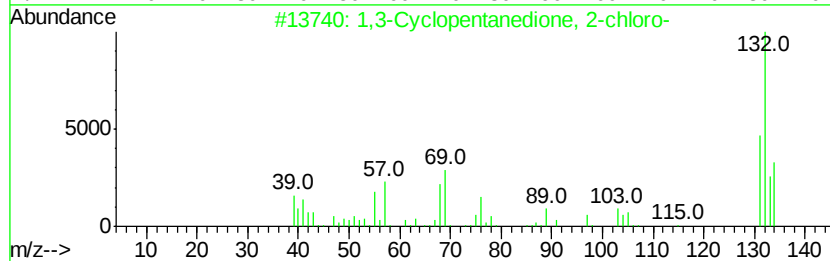
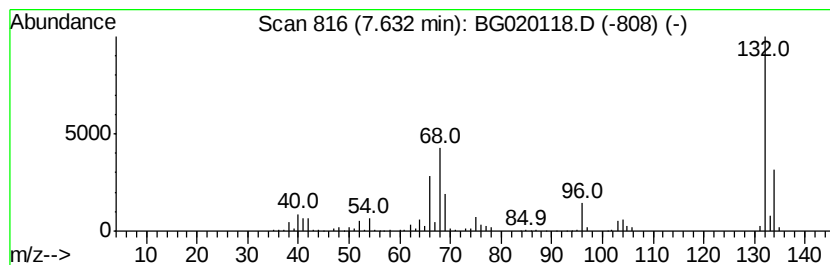
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown7.63 Concentration Rank 1

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.63 | 76.40 ng | 436044 | 1,4-Dichlorobenzene-d4 | 8.10 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|---|----------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1,3-Cyclopentanedione, 2-chloro- | 132 | C5H5ClO2  | 014203-19-1  | 35   |
| 2    |    |   | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 22   |
| 3    |    |   | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 16   |
| 4    |    |   | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 12   |
| 5    |    |   | Benzene, 1,3,5-trifluoro-        | 132 | C6H3F3    | 000372-38-3  | 10   |



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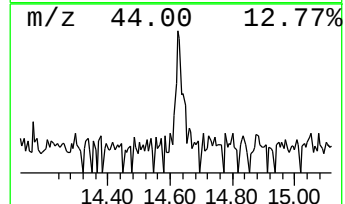
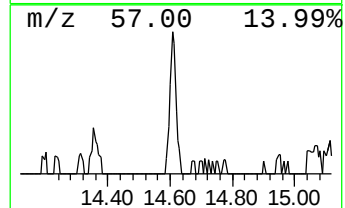
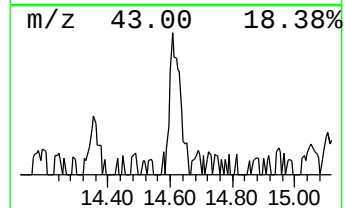
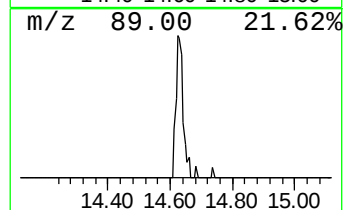
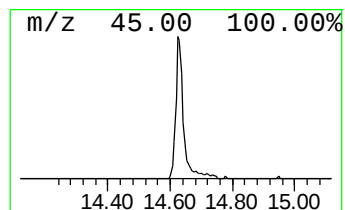
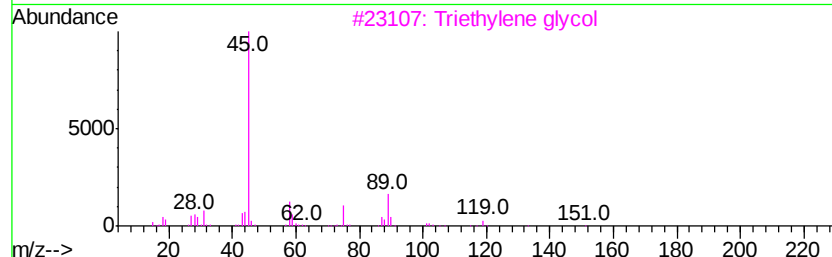
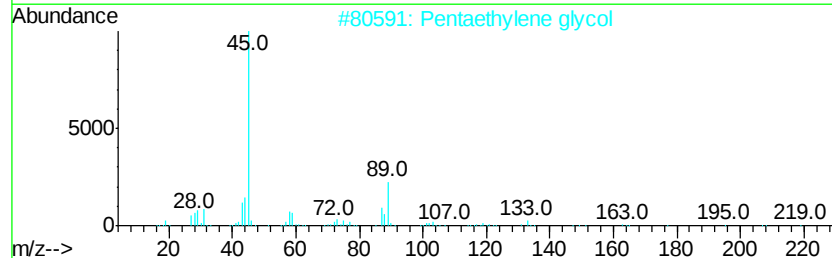
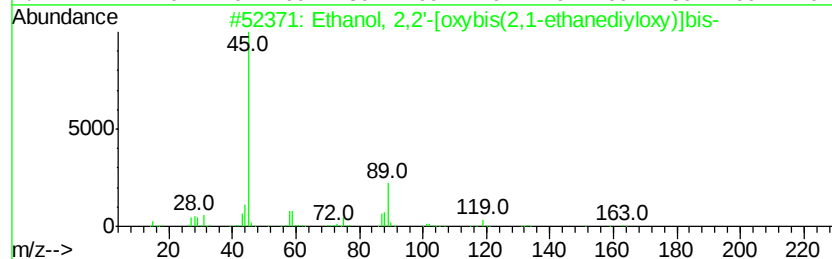
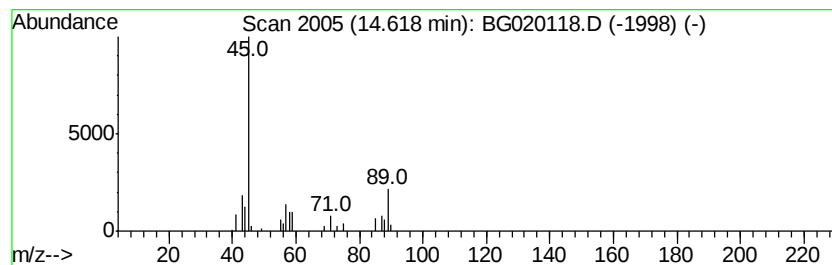
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Peak Number 6 Ethanol, 2,2'-[oxybis(2,1-e... Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.62 | 2.75 ng | 38224 | Acenaphthene-d10 | 14.73 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Ethanol, 2,2'-[oxybis(2,1-ethane... | 194 | C8H18O5  | 000112-60-7 | 83   |
| 2    |    | Pentaethylene glycol                | 238 | C10H22O6 | 004792-15-8 | 74   |
| 3    |    | Triethylene glycol                  | 150 | C6H14O4  | 000112-27-6 | 64   |
| 4    |    | Ethanol, 2-[2-(2-methoxyethoxy)e... | 164 | C7H16O4  | 000112-35-6 | 40   |
| 5    |    | Ethanol, 2-[2-(2-butoxyethoxy)et... | 206 | C10H22O4 | 000143-22-6 | 40   |



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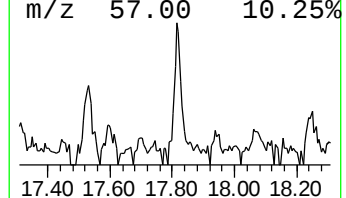
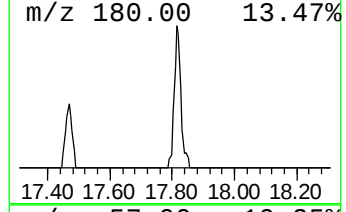
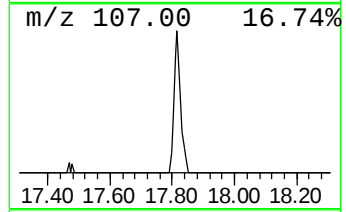
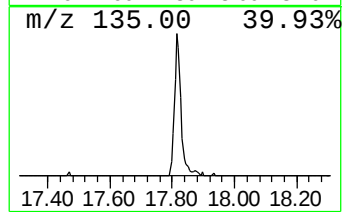
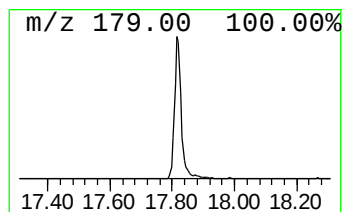
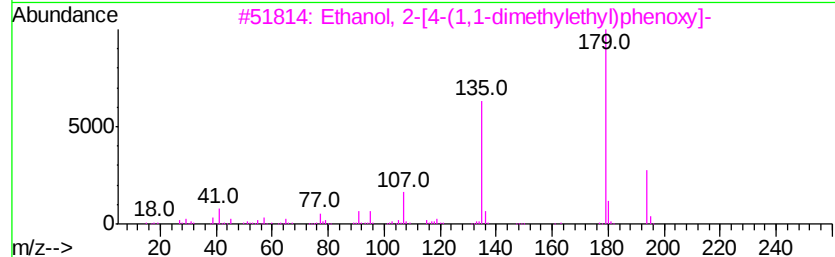
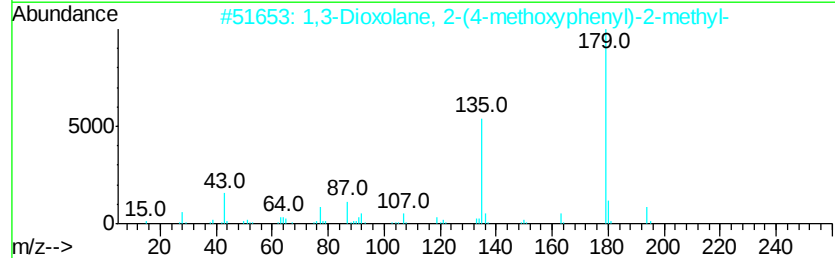
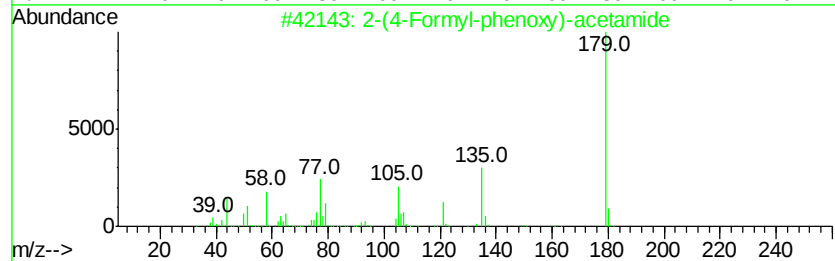
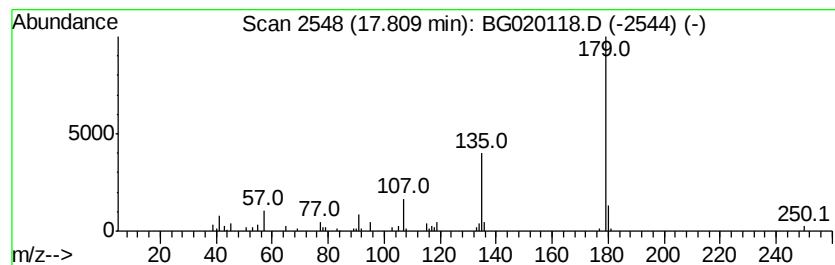
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 2-(4-Formyl-phenoxy)-acetamide Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 17.81 | 3.88 ng | 73576 | Phenanthrene-d10 | 17.47 |

| Hit# | of | Tentative ID                          | MW  | MolForm   | CAS#        | Qual |
|------|----|---------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 2-(4-Formyl-phenoxy)-acetamide        | 179 | C9H9NO3   | 135857-20-4 | 64   |
| 2    |    | 1,3-Dioxolane, 2-(4-methoxypheny...   | 194 | C11H14O3  | 036881-00-2 | 64   |
| 3    |    | Ethanol, 2-[4-(1,1-dimethylethyl)...] | 194 | C12H18O2  | 000713-46-2 | 50   |
| 4    |    | Benzothiazole-6-carboxylic acid       | 179 | C8H5NO2S  | 003622-35-3 | 40   |
| 5    |    | Acetic acid, [2-[(2-propenylamin...   | 235 | C12H13NO4 | 000119-45-9 | 37   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG121215\  
Data File : BG020118.D  
Acq On : 11 Dec 2015 22:43  
Operator : UM/SJ  
Sample : G4729-15  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
W-42

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

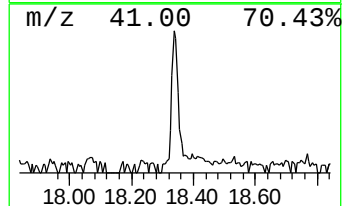
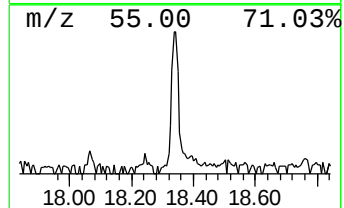
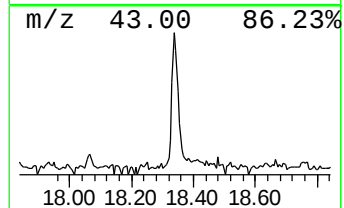
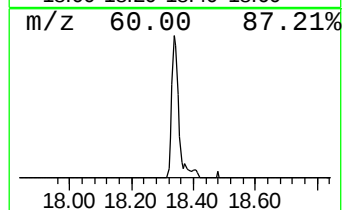
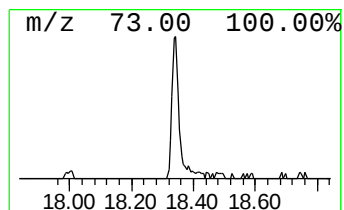
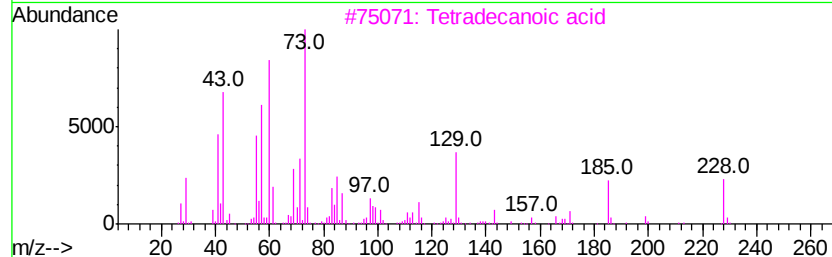
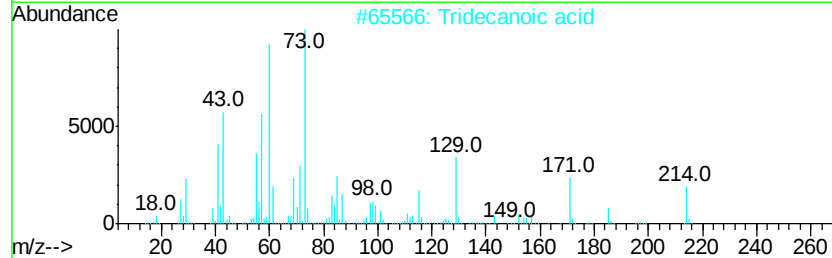
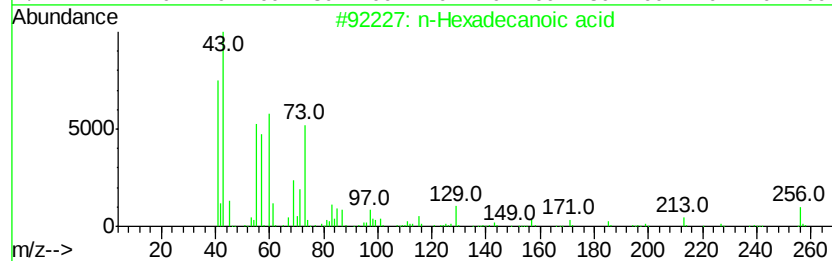
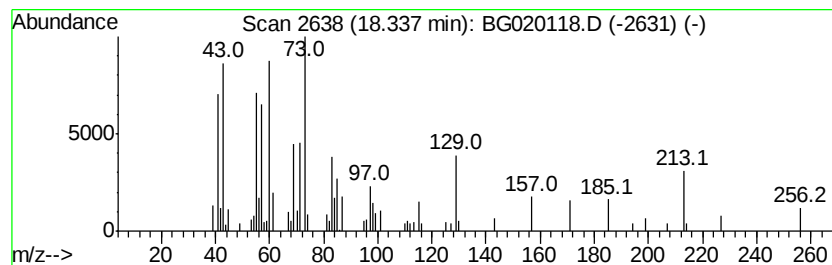
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 n-Hexadecanoic acid Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 18.34 | 3.85 ng | 72946 | Phenanthrene-d10 | 17.47 |

| Hit# of 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|-----------|---------------------|-----|----------|-------------|------|
| 1         | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 95   |
| 2         | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 64   |
| 3         | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 64   |
| 4         | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 50   |
| 5         | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 49   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG121215\  
Data File : BG020118.D  
Acq On : 11 Dec 2015 22:43  
Operator : UM/SJ  
Sample : G4729-15  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

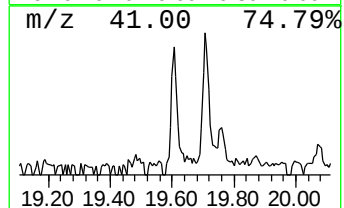
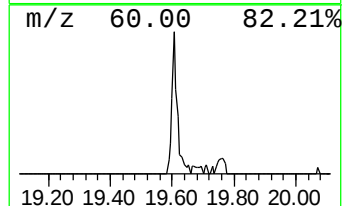
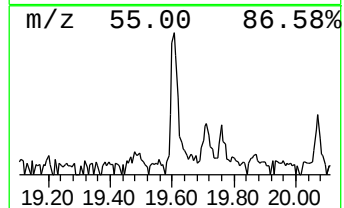
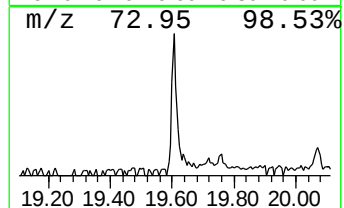
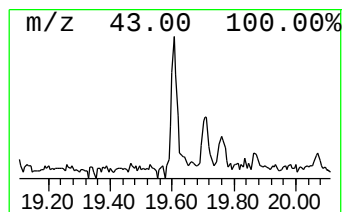
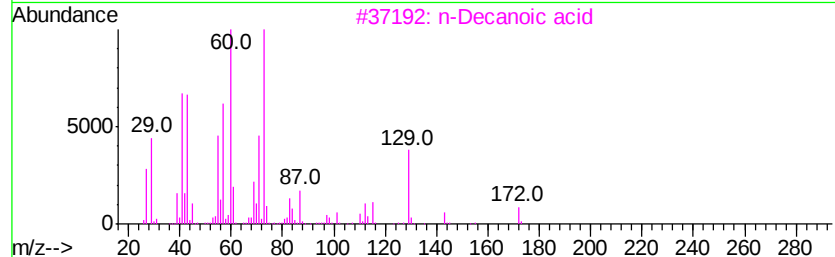
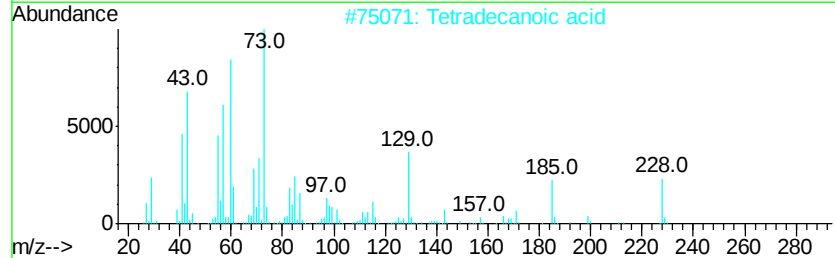
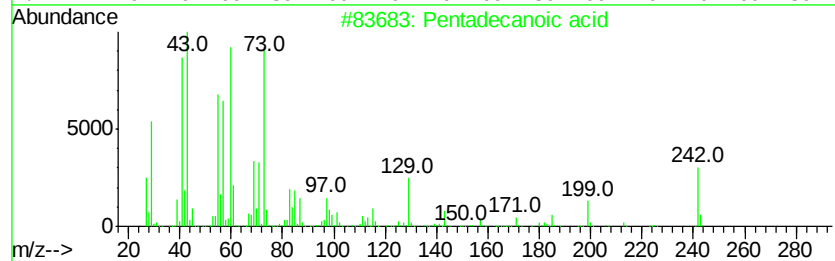
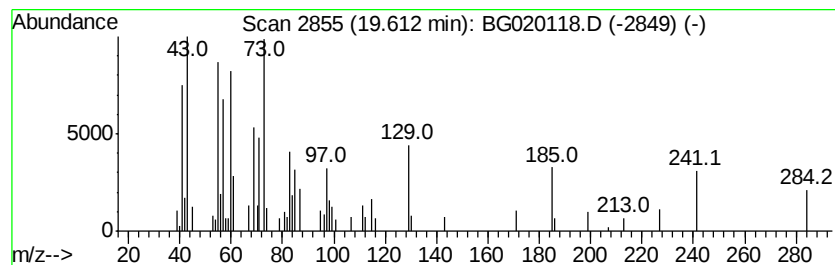
Instrument :  
BNA\_G  
ClientSampleId :  
W-42

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.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 9 Pentadecanoic acid Concentration Rank 11

| R.T.      | EstConc            | Area  | Relative to ISTD |             | R.T.  |
|-----------|--------------------|-------|------------------|-------------|-------|
| 19.61     | 2.45 ng            | 57209 | Chrysene-d12     |             | 21.74 |
| Hit# of 5 | Tentative ID       | MW    | MolForm          | CAS#        | Qual  |
| 1         | Pentadecanoic acid | 242   | C15H30O2         | 001002-84-2 | 80    |
| 2         | Tetradecanoic acid | 228   | C14H28O2         | 000544-63-8 | 72    |
| 3         | n-Decanoic acid    | 172   | C10H20O2         | 000334-48-5 | 64    |
| 4         | Tridecanoic acid   | 214   | C13H26O2         | 000638-53-9 | 53    |
| 5         | Undecanoic acid    | 186   | C11H22O2         | 000112-37-8 | 49    |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG121215\  
Data File : BG020118.D  
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Operator : UM/SJ  
Sample : G4729-15  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
W-42

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

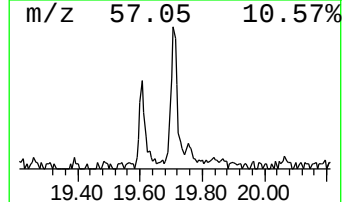
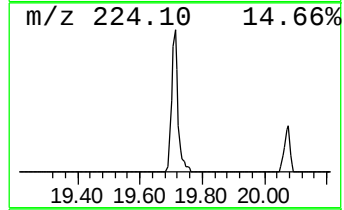
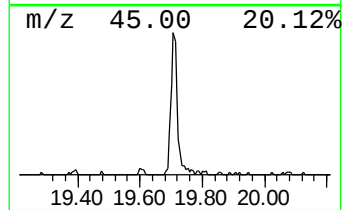
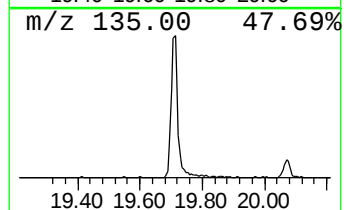
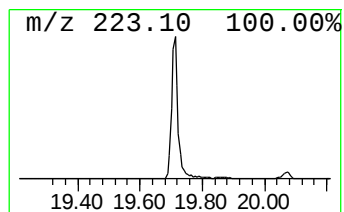
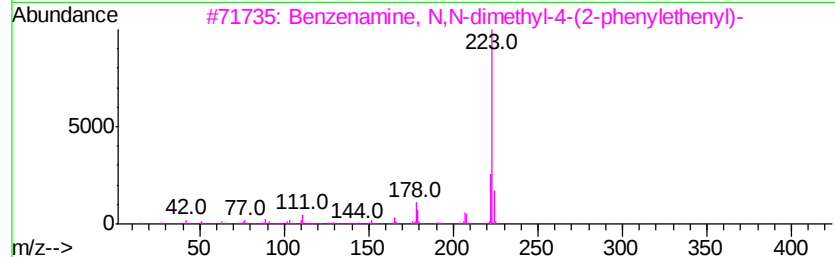
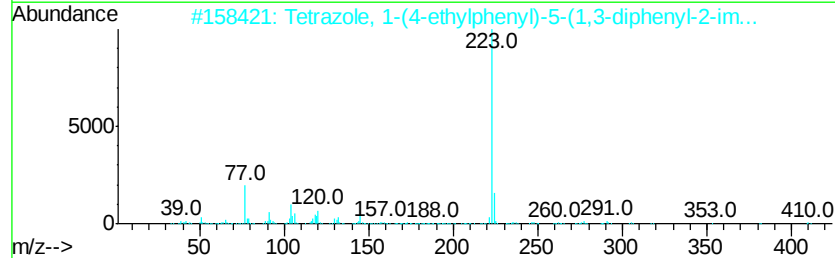
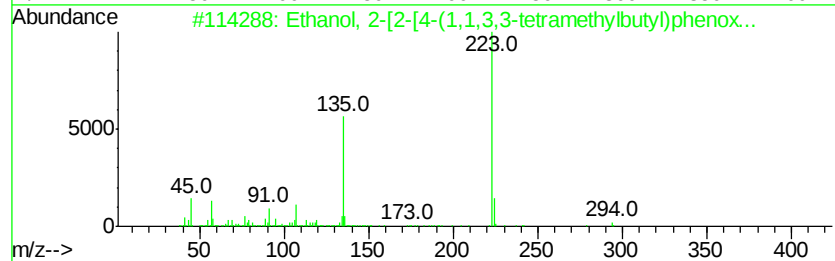
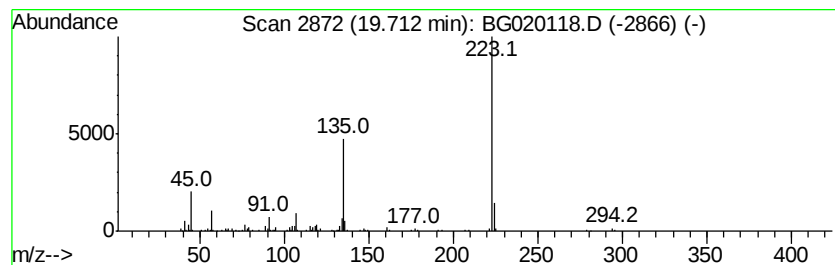
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.71 | 7.39 ng | 172811 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Ethanol, 2-[2-[4-(1,1,3,3-tetram... | 294 | C18H30O3 | 002315-61-9 | 96   |
| 2    |    | Tetrazole, 1-(4-ethylphenyl)-5-(... | 410 | C25H26N6 | 125038-06-4 | 38   |
| 3    |    | Benzenamine, N,N-dimethyl-4-(2-p... | 223 | C16H17N  | 001145-73-9 | 38   |
| 4    |    | Benzenamine, N,N-dimethyl-4-(2-p... | 223 | C16H17N  | 000838-95-9 | 32   |
| 5    |    | 4,4'-Trimethylenebis(1-methylpip... | 238 | C15H30N2 | 064168-11-2 | 32   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG121215\  
Data File : BG020118.D  
Acq On : 11 Dec 2015 22:43  
Operator : UM/SJ  
Sample : G4729-15  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
W-42

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

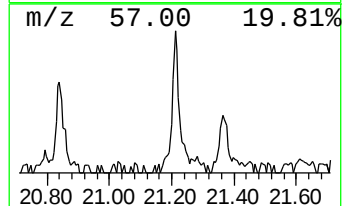
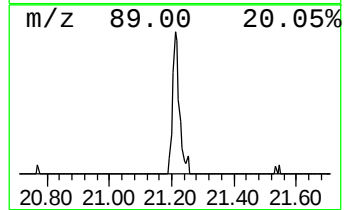
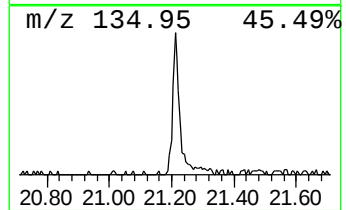
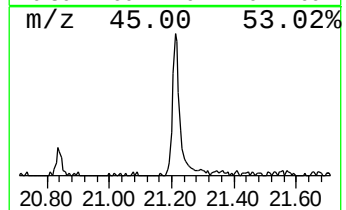
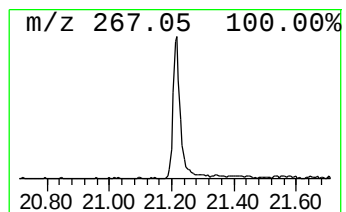
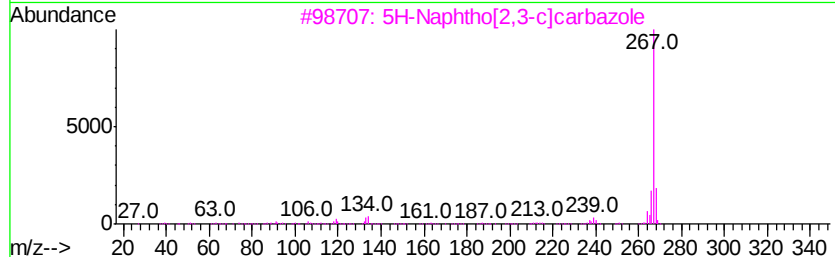
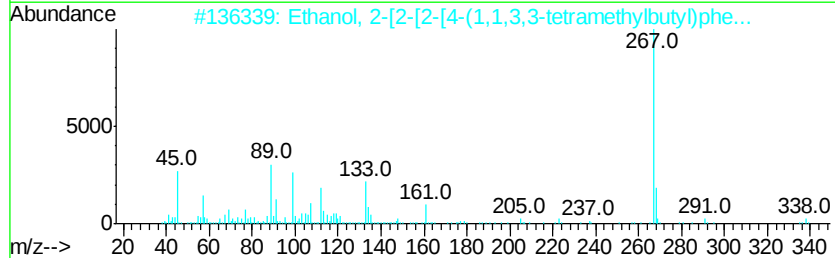
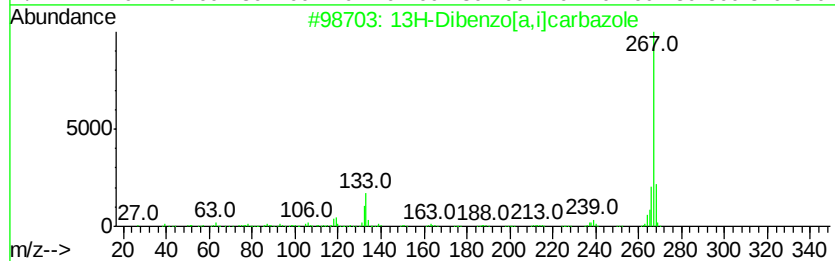
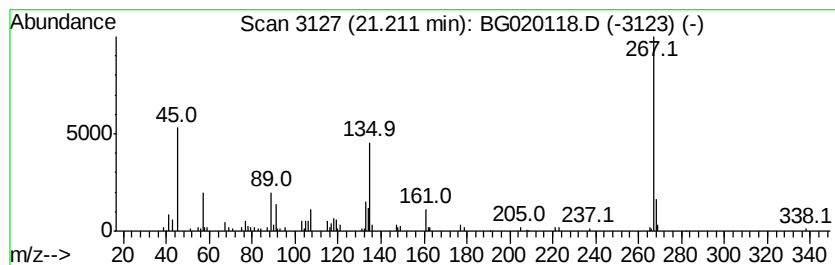
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 unknown21.21 Concentration Rank 7

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 21.21 | 3.97 ng | 92806 | Chrysene-d12     | 21.74 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 13H-Dibenzo[a,i]carbazole           | 267 | C20H13N  | 000239-64-5 | 46   |
| 2    |    | Ethanol, 2-[2-[2-[4-(1,1,3,3-tet... | 338 | C20H34O4 | 002315-62-0 | 45   |
| 3    |    | 5H-Naphtho[2,3-c]carbazole          | 267 | C20H13N  | 000198-96-9 | 35   |
| 4    |    | 7H-Dibenzo[c,q]carbazole            | 267 | C20H13N  | 000194-59-2 | 27   |
| 5    |    | 1,5-Diphenyl-3-methylthio-1,2,4-... | 267 | C15H13NS | 072234-23-2 | 25   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121215\  
Data File : BG020118.D  
Acq On : 11 Dec 2015 22:43  
Operator : UM/SJ  
Sample : G4729-15  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
W-42

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG120215.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 |
| Butane, 2-methoxy... | 3.20  | 9.3     | ng    | 53346    | 1                     | 8.10  | 114148 | 20.0 |
| unknown5.18          | 5.18  | 8.8     | ng    | 50003    | 1                     | 8.10  | 114148 | 20.0 |
| unknown7.63          | 7.63  | 76.4    | ng    | 436044   | 1                     | 8.10  | 114148 | 20.0 |
| Ethanol, 2,2'-[ox... | 14.62 | 2.8     | ng    | 38224    | 3                     | 14.73 | 278030 | 20.0 |
| 2-(4-Formyl-pheno... | 17.81 | 3.9     | ng    | 73576    | 4                     | 17.47 | 379020 | 20.0 |
| n-Hexadecanoic acid  | 18.34 | 3.9     | ng    | 72946    | 4                     | 17.47 | 379020 | 20.0 |
| Pentadecanoic acid   | 19.61 | 2.5     | ng    | 57209    | 5                     | 21.74 | 467861 | 20.0 |
| Ethanol, 2-[2-[4-... | 19.71 | 7.4     | ng    | 172811   | 5                     | 21.74 | 467861 | 20.0 |
| unknown21.21         | 21.21 | 4.0     | ng    | 92806    | 5                     | 21.74 | 467861 | 20.0 |