

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
 Data File : BG032589.D  
 Acq On : 7 Feb 2018 8:04  
 Operator : SJ/JU  
 Sample : J1455-09  
 Misc : MDL-W-8 ppm  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 LW-04-B

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-BG012918.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         | 3.367       | 6             | 13          | 22           | rBV5     | 16205          | 38409         | 1.71%           | 0.416%        |
| 2         | 3.461       | 23            | 29          | 40           | rVB      | 59660          | 99480         | 4.44%           | 1.078%        |
| 3         | 6.146       | 478           | 486         | 499          | rBV      | 101571         | 199000        | 8.88%           | 2.157%        |
| 4         | 7.826       | 766           | 772         | 789          | rBV2     | 52994          | 128557        | 5.74%           | 1.393%        |
| 5         | 8.220       | 832           | 839         | 854          | rBV      | 251523         | 500445        | 22.34%          | 5.424%        |
| 6         | 8.708       | 915           | 922         | 931          | rBV      | 45514          | 97198         | 4.34%           | 1.053%        |
| 7         | 9.043       | 971           | 979         | 996          | rBV      | 233129         | 486528        | 21.72%          | 5.273%        |
| 8         | 9.906       | 1118          | 1126        | 1140         | rBV2     | 188346         | 399194        | 17.82%          | 4.327%        |
| 9         | 11.581      | 1405          | 1411        | 1422         | rBV      | 65313          | 140970        | 6.29%           | 1.528%        |
| 10        | 13.966      | 1810          | 1817        | 1832         | rBV      | 792845         | 1324351       | 59.12%          | 14.354%       |
| 11        | 15.341      | 2044          | 2051        | 2057         | rBV2     | 171866         | 293821        | 13.12%          | 3.185%        |
| 12        | 15.494      | 2073          | 2077        | 2082         | rBV6     | 22478          | 39081         | 1.74%           | 0.424%        |
| 13        | 15.823      | 2126          | 2133        | 2139         | rBV4     | 26724          | 53852         | 2.40%           | 0.584%        |
| 14        | 15.940      | 2148          | 2153        | 2157         | rVB8     | 13476          | 22513         | 1.01%           | 0.244%        |
| 15        | 16.117      | 2179          | 2183        | 2187         | rBV7     | 13799          | 27165         | 1.21%           | 0.294%        |
| 16        | 16.164      | 2187          | 2191        | 2193         | rVV5     | 23050          | 40405         | 1.80%           | 0.438%        |
| 17        | 16.211      | 2195          | 2199        | 2210         | rVB3     | 38830          | 91511         | 4.09%           | 0.992%        |
| 18        | 16.687      | 2273          | 2280        | 2287         | rBV3     | 120945         | 277273        | 12.38%          | 3.005%        |
| 19        | 16.816      | 2296          | 2302        | 2314         | rBV      | 693279         | 1202463       | 53.68%          | 13.033%       |
| 20        | 18.073      | 2510          | 2516        | 2526         | rVB      | 226957         | 384315        | 17.16%          | 4.165%        |
| 21        | 18.890      | 2652          | 2655        | 2661         | rBV5     | 29160          | 50592         | 2.26%           | 0.548%        |
| 22        | 20.611      | 2939          | 2948        | 2955         | rBV      | 1681497        | 2239990       | 100.00%         | 24.278%       |
| 23        | 21.945      | 3170          | 3175        | 3182         | rBV9     | 21071          | 42726         | 1.91%           | 0.463%        |
| 24        | 22.397      | 3245          | 3252        | 3266         | rVB2     | 261370         | 518868        | 23.16%          | 5.624%        |
| 25        | 26.070      | 3863          | 3877        | 3891         | rBV2     | 162333         | 527646        | 23.56%          | 5.719%        |

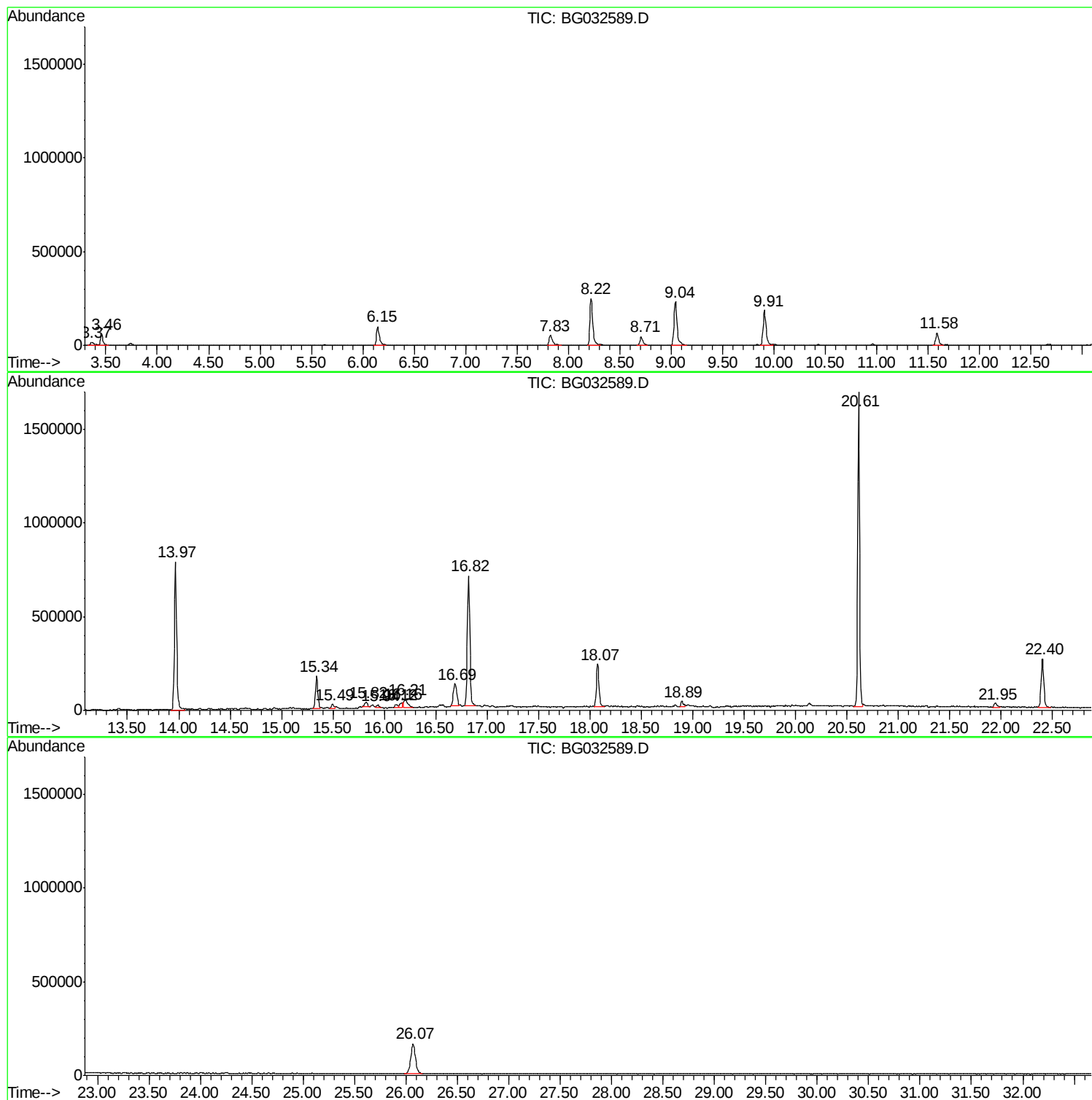
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LW-04-B

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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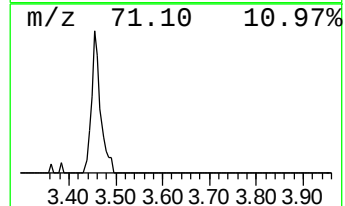
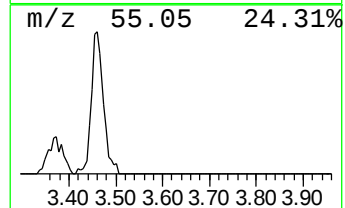
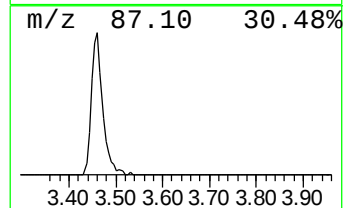
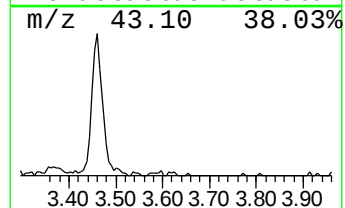
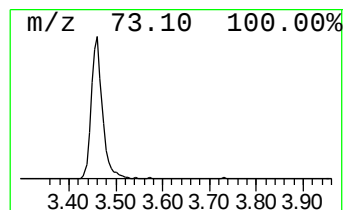
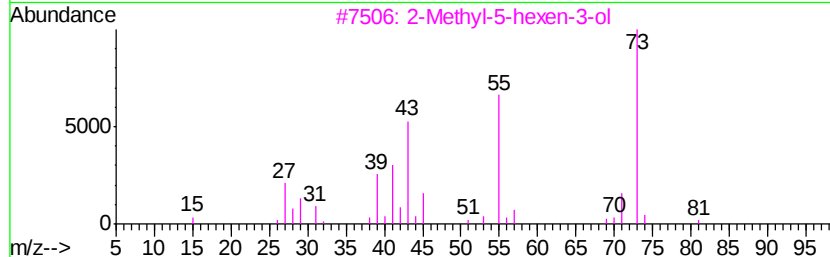
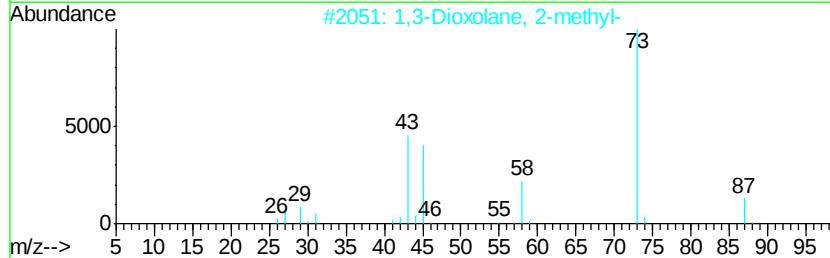
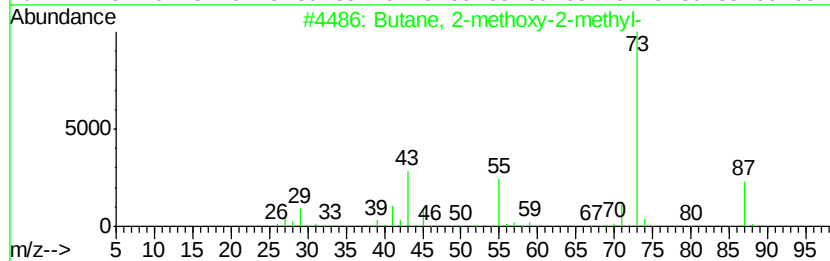
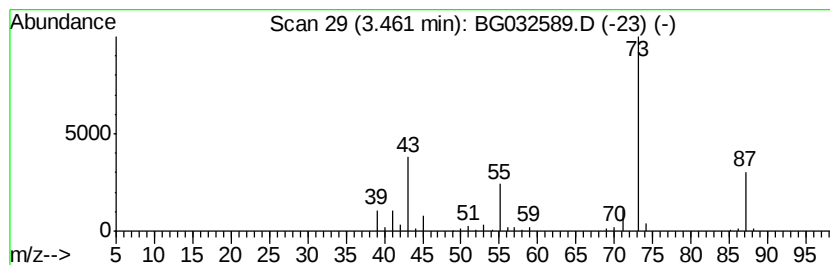
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T. | EstConc  | Area  | Relative to ISTD       | R.T. |
|------|----------|-------|------------------------|------|
| 3.46 | 20.47 ng | 99480 | 1,4-Dichlorobenzene-d4 | 8.70 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 2    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 23   |
| 3    |    | 2-Methyl-5-hexen-3-ol       | 114 | C7H14O  | 032815-70-6 | 12   |
| 4    |    | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 5    |    | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 9    |



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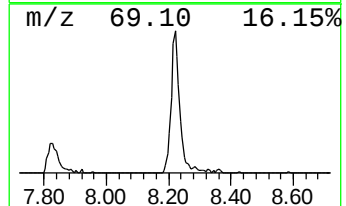
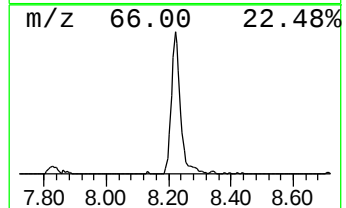
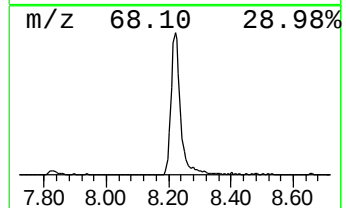
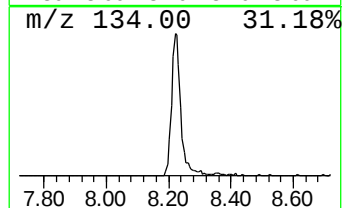
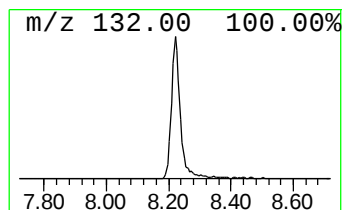
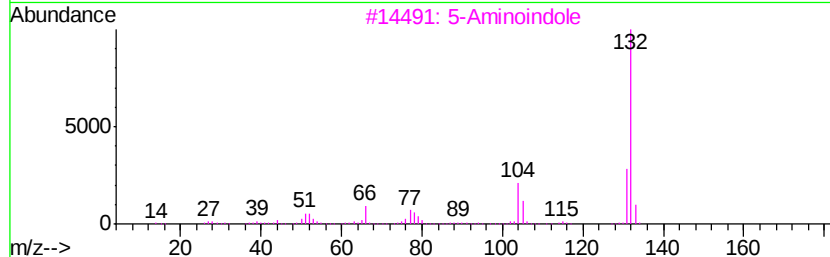
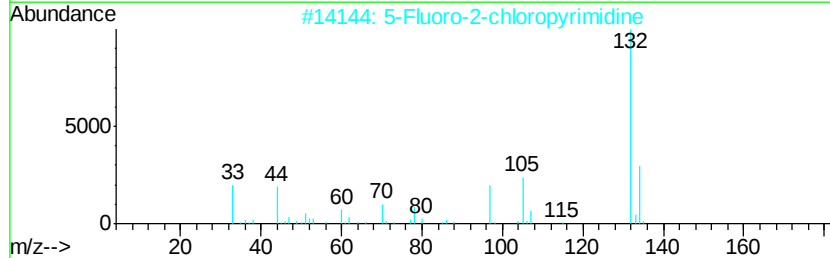
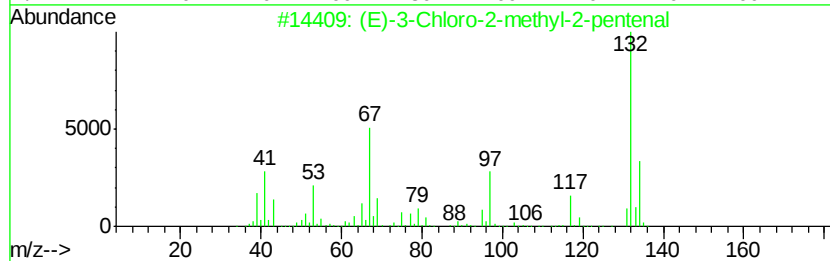
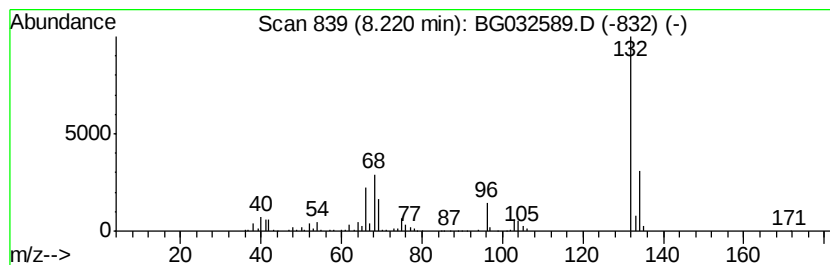
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Peak Number 3 unknown8.22 Concentration Rank 1

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 8.22 | 102.97 ng | 500445 | 1,4-Dichlorobenzene-d4 | 8.70 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 25   |
| 2    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 17   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 16   |
| 4    |    | (5-Methyl-2-pyridyl)acetonitrile | 132 | C8H8N2    | 1000241-93-9 | 12   |
| 5    |    | 4-Chloro-2-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-00-5  | 9    |



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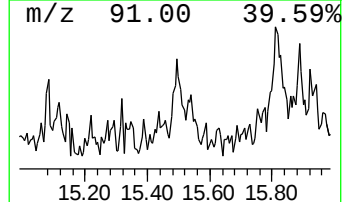
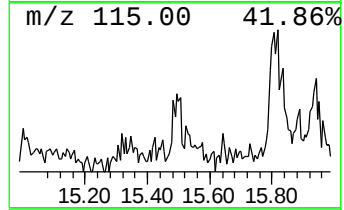
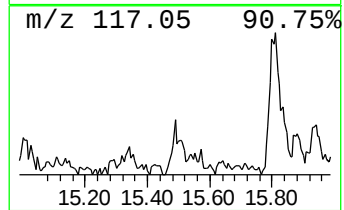
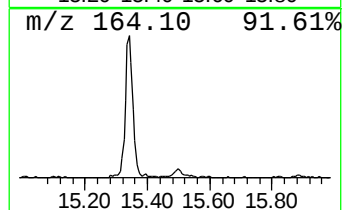
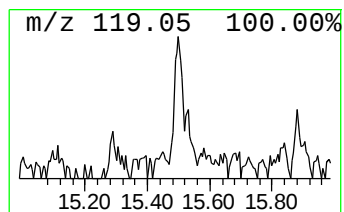
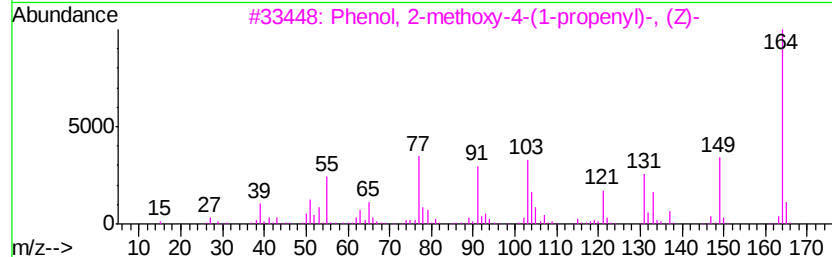
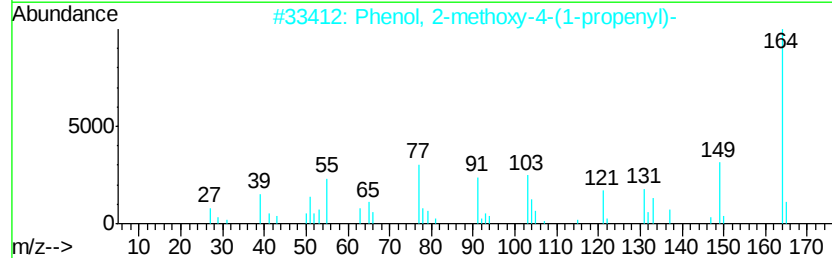
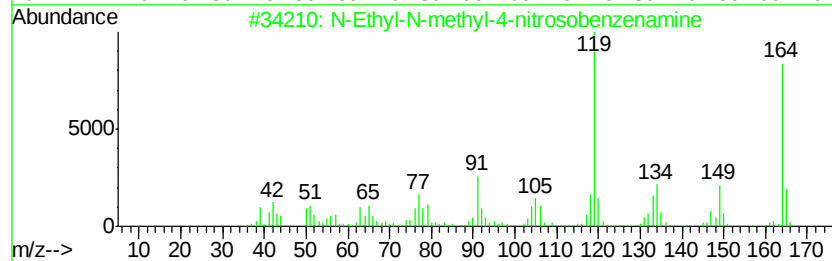
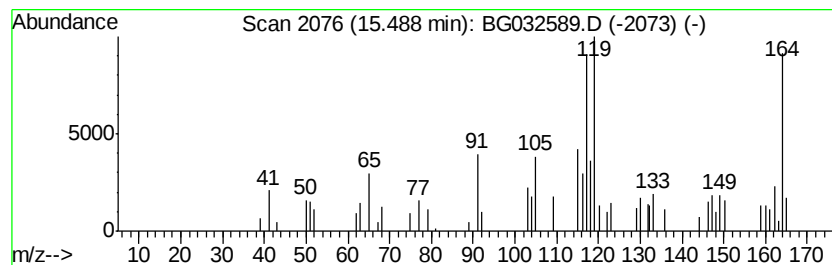
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 N-Ethyl-N-methyl-4-nitrosob... Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.49 | 2.66 ng | 39081 | Acenaphthene-d10 | 15.34 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|---------|---|-------------------------------------|-----|----------|-------------|------|
| 1       |   | N-Ethyl-N-methyl-4-nitrosobenzen... | 164 | C9H12N2O | 036479-98-8 | 64   |
| 2       |   | Phenol, 2-methoxy-4-(1-propenyl)-   | 164 | C10H12O2 | 000097-54-1 | 25   |
| 3       |   | Phenol, 2-methoxy-4-(1-propenyl)... | 164 | C10H12O2 | 005912-86-7 | 25   |
| 4       |   | Pvridine, 5-ethenyl-2-methyl-       | 119 | C8H9N    | 000140-76-1 | 25   |
| 5       |   | Eugenol                             | 164 | C10H12O2 | 000097-53-0 | 25   |



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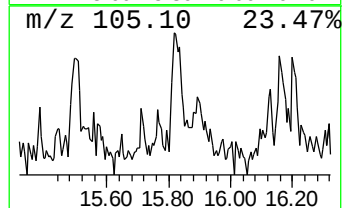
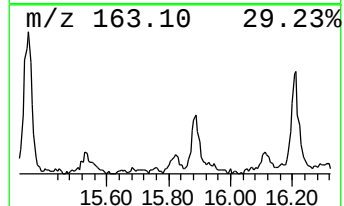
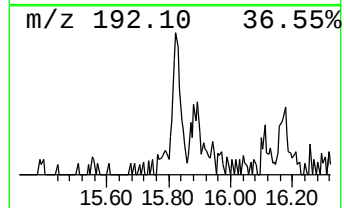
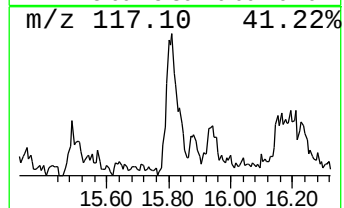
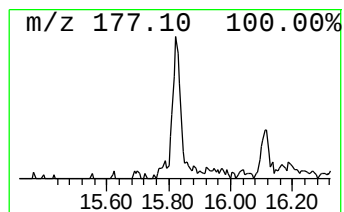
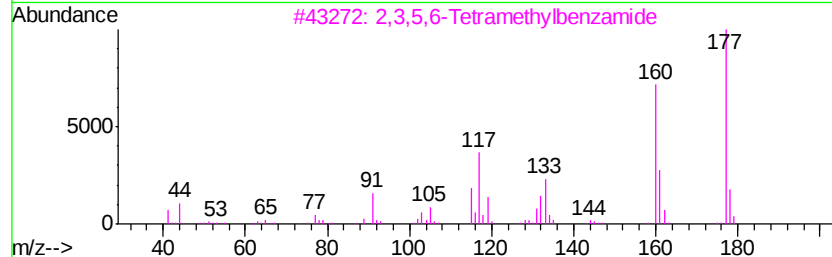
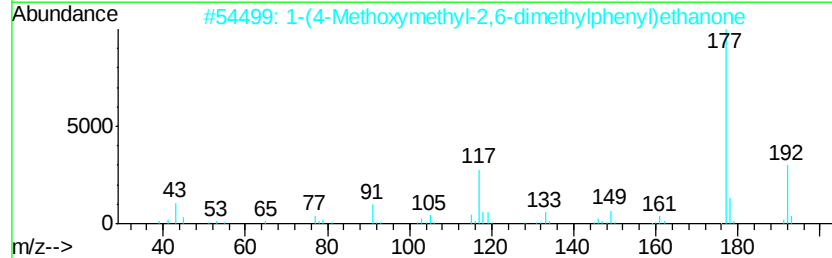
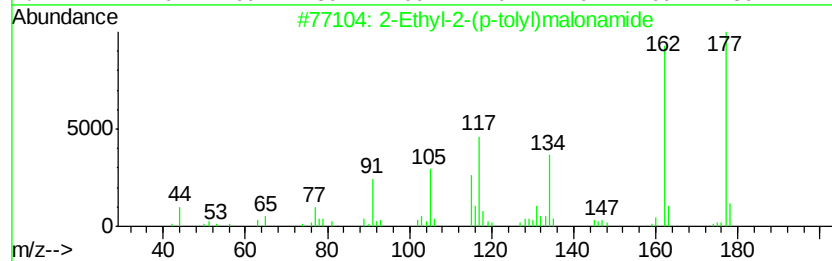
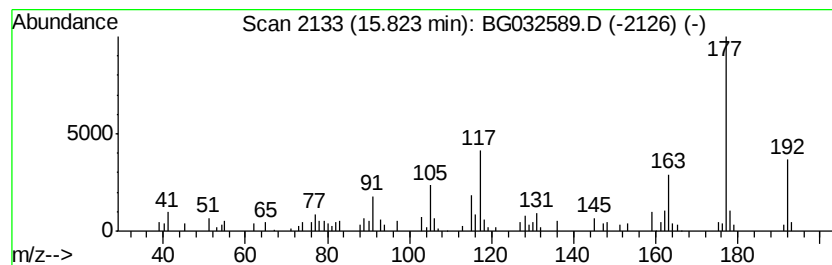
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\*\*\*\*\*  
Peak Number 6 2-Ethyl-2-(p-tolyl)malonamide Concentration Rank 7

| R.T.    | EstConc                             | Area         | Relative to ISTD |              | R.T.  |      |
|---------|-------------------------------------|--------------|------------------|--------------|-------|------|
| 15.82   | 3.67 ng                             | 53852        | Acenaphthene-d10 |              | 15.34 |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm      | CAS#  | Qual |
| 1       | 2-Ethyl-2-(p-tolyl)malonamide       | 220          | C12H16N2O2       | 068692-83-1  | 53    |      |
| 2       | 1-(4-Methoxymethyl-2,6-dimethylp... | 192          | C12H16O2         | 1000202-02-2 | 49    |      |
| 3       | 2,3,5,6-Tetramethylbenzamide        | 177          | C11H15NO         | 099858-56-7  | 49    |      |
| 4       | 2,6-Diisopropylanisole              | 192          | C13H20O          | 002944-52-7  | 46    |      |
| 5       | Pyrazine, 2,5-bis(1,1-dimethylet... | 192          | C12H20N2         | 018709-51-8  | 46    |      |



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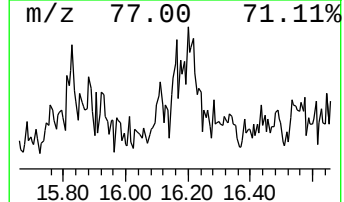
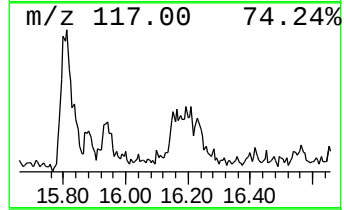
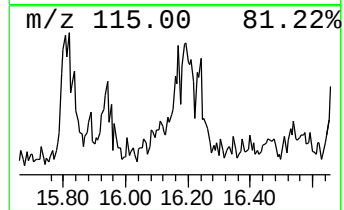
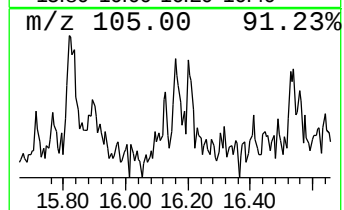
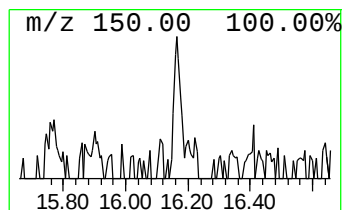
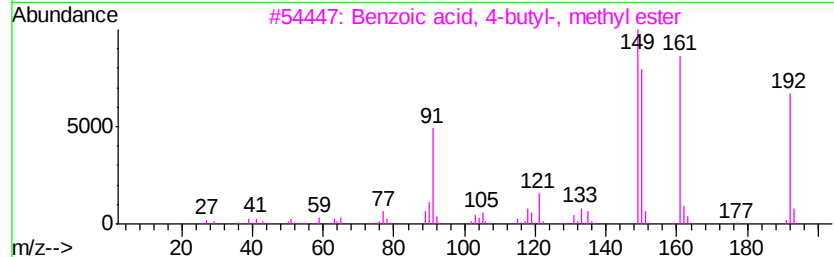
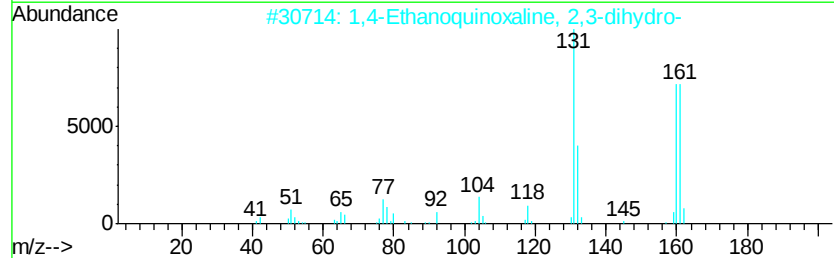
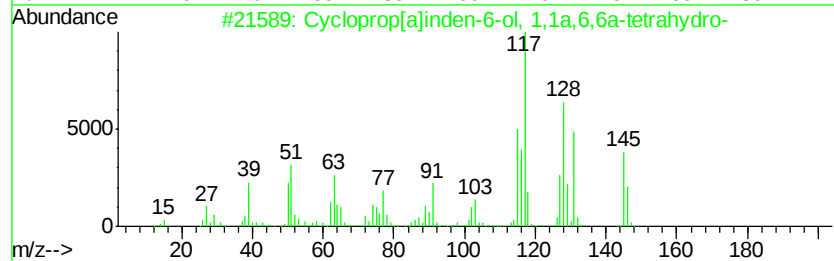
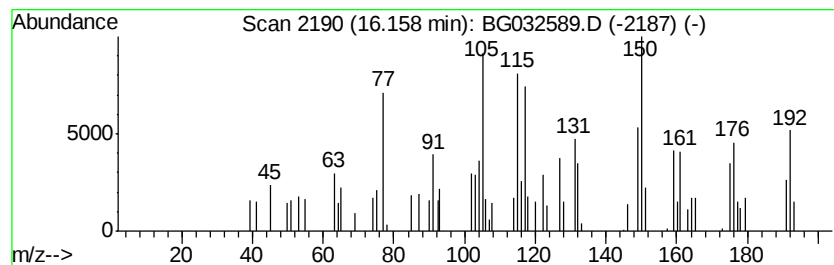
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 unknown16.16 Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.16 | 2.75 ng | 40405 | Acenaphthene-d10 | 15.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Cycloprop[alinden-6-ol, 1,1a,6,6... | 146 | C10H10O  | 022228-27-9 | 30   |
| 2    |    | 1,4-Ethanoquinoxaline, 2,3-dihydro- | 160 | C10H12N2 | 007140-45-6 | 14   |
| 3    |    | Benzoic acid, 4-butyl-, methyl e... | 192 | C12H16O2 | 020651-69-8 | 14   |
| 4    |    | Thiophene, tetrahydro-3-phenyl-,... | 180 | C10H12OS | 093134-22-6 | 10   |
| 5    |    | 3a,6-Methano-3aH-indene, 2,3,6,7... | 132 | C10H12   | 098640-29-0 | 10   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032589.D  
Acq On : 7 Feb 2018 8:04  
Operator : SJ/JU  
Sample : J1455-09  
Misc : MDL-W-8 ppm  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-04-B

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

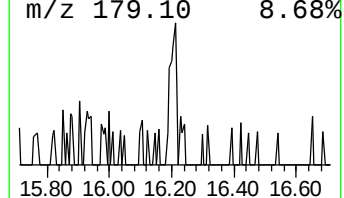
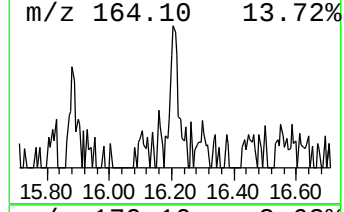
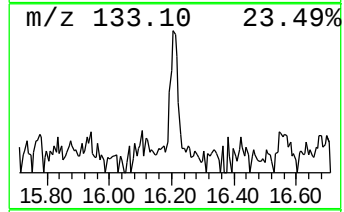
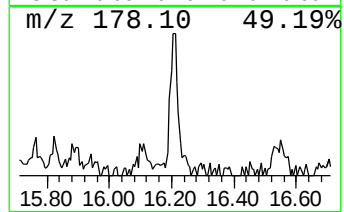
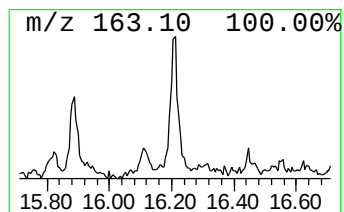
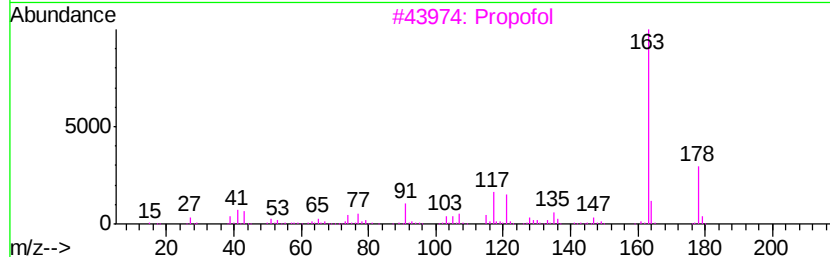
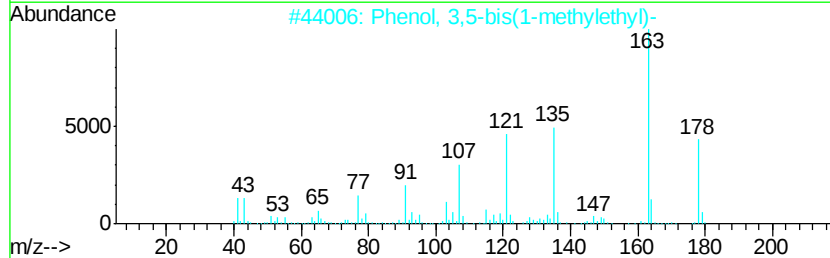
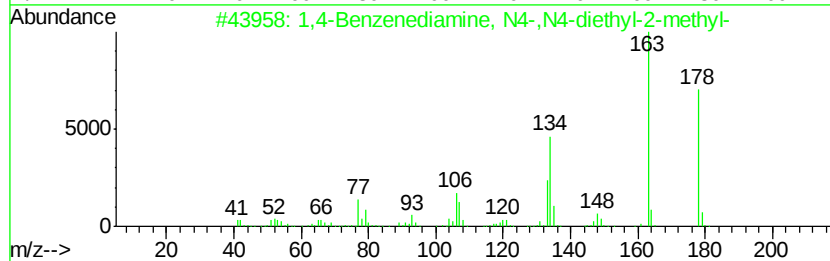
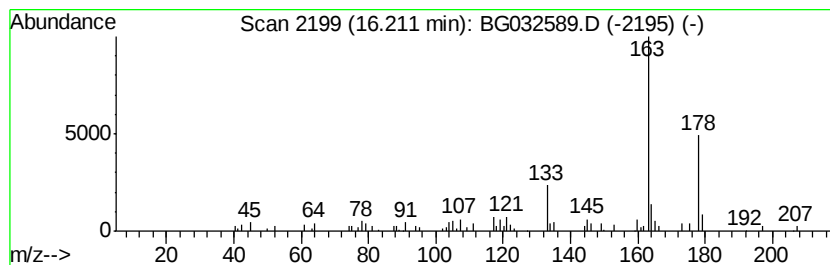
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 1,4-Benzenediamine, N4-,N4-... Concentration Rank 6

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.21 | 6.23 ng | 91511 | Acenaphthene-d10 | 15.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,4-Benzenediamine, N4-,N4-dieth... | 178 | C11H18N2 | 000148-71-0 | 72   |
| 2    |    | Phenol, 3,5-bis(1-methylethyl)-     | 178 | C12H18O  | 026886-05-5 | 64   |
| 3    |    | Propofol                            | 178 | C12H18O  | 002078-54-8 | 64   |
| 4    |    | 6-tert-Butyl-2,4-dimethylphenol     | 178 | C12H18O  | 001879-09-0 | 64   |
| 5    |    | Ethanone, 1-(2,3-dihydro-1,4-ben... | 178 | C10H10O3 | 002879-20-1 | 59   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032589.D  
Acq On : 7 Feb 2018 8:04  
Operator : SJ/JU  
Sample : J1455-09  
Misc : MDL-W-8 ppm  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-04-B

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

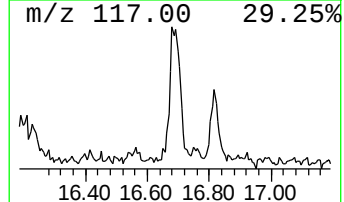
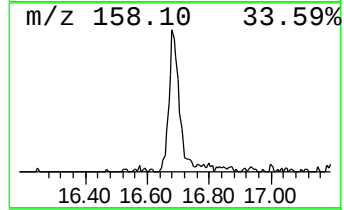
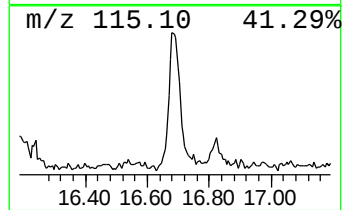
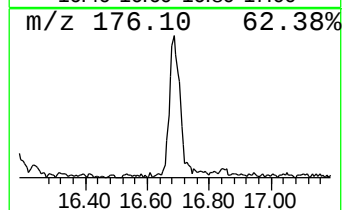
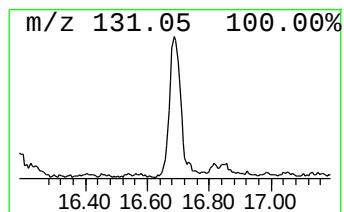
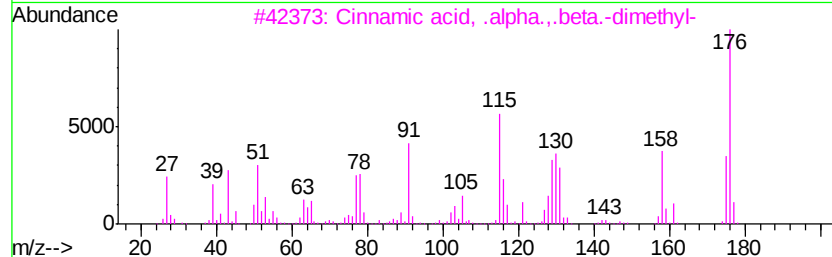
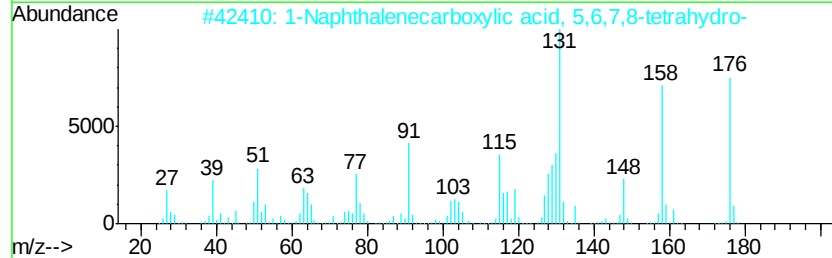
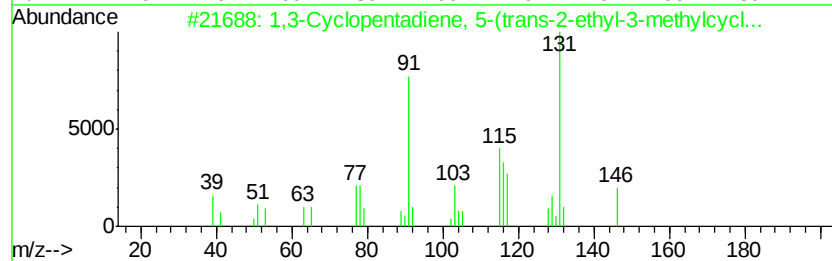
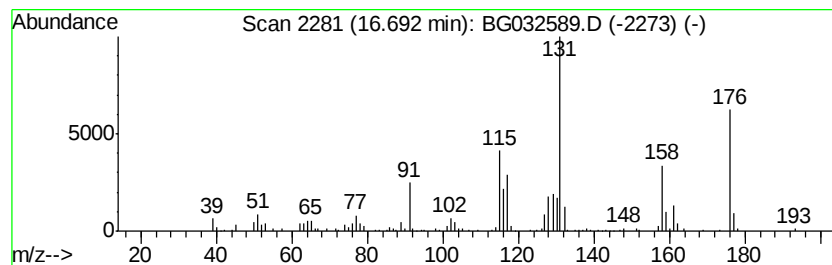
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown16.69 Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 16.69 | 18.87 ng | 277273 | Acenaphthene-d10 | 15.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 1,3-Cyclopentadiene, 5-(trans-2-... | 146 | C11H14   | 079209-36-2  | 47   |
| 2    |    | 1-Naphthalenecarboxylic acid, 5,... | 176 | C11H12O2 | 004242-18-6  | 47   |
| 3    |    | Cinnamic acid, .alpha.,.beta.-di... | 176 | C11H12O2 | 1000151-56-4 | 43   |
| 4    |    | 3-Pentenoic acid, 4-phenyl-         | 176 | C11H12O2 | 053774-19-9  | 43   |
| 5    |    | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14   | 053172-84-2  | 37   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG020618\  
Data File : BG032589.D  
Acq On : 7 Feb 2018 8:04  
Operator : SJ/JU  
Sample : J1455-09  
Misc : MDL-W-8 ppm  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-04-B

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

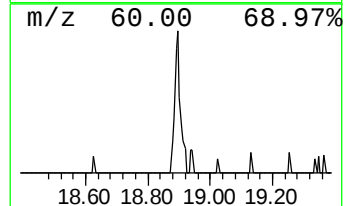
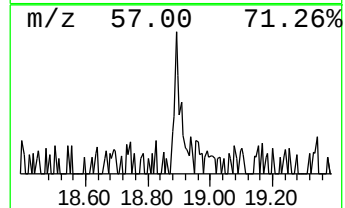
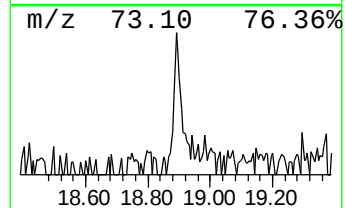
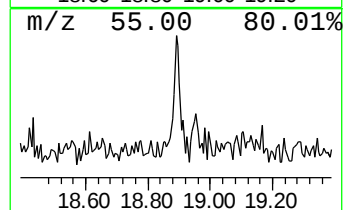
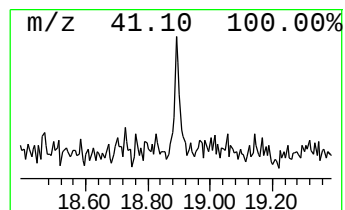
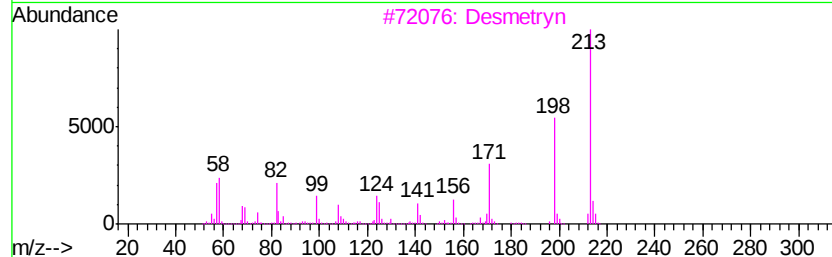
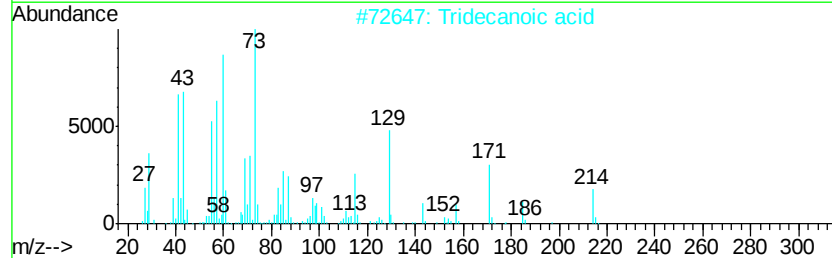
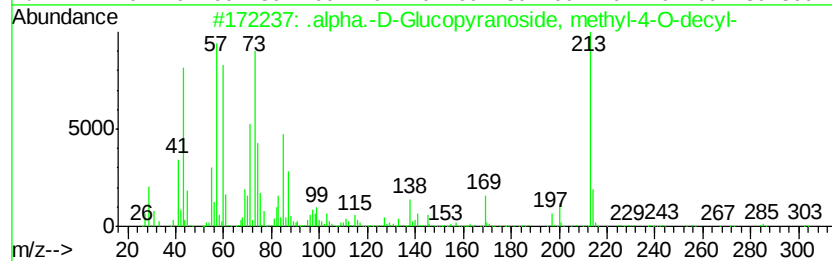
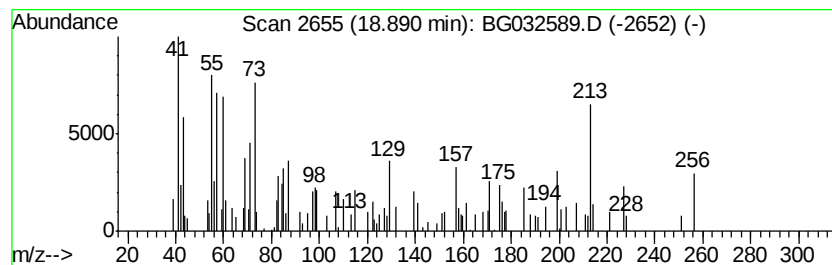
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 unknown18.89 Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 18.89 | 2.63 ng | 50592 | Phenanthrene-d10 | 18.07 |

| Hit# | of | 5                                   | Tentative ID | MW           | MolForm      | CAS# | Qual |
|------|----|-------------------------------------|--------------|--------------|--------------|------|------|
| 1    | .  | alpha.-D-Glucopyranoside, methy...  | 334          | C17H34O6     | 1000157-24-7 | 30   |      |
| 2    |    | Tridecanoic acid                    | 214          | C13H26O2     | 000638-53-9  | 30   |      |
| 3    |    | Desmetryn                           | 213          | C8H15N5S     | 001014-69-3  | 25   |      |
| 4    |    | n-Hexadecanoic acid                 | 256          | C16H32O2     | 000057-10-3  | 22   |      |
| 5    |    | Benzenamine, 2-chloro-N,N-diethy... | 228          | C10H13ClN2O2 | 000086-49-7  | 18   |      |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG020618\  
Data File : BG032589.D  
Acq On : 7 Feb 2018 8:04  
Operator : SJ/JU  
Sample : J1455-09  
Misc : MDL-W-8 ppm  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
LW-04-B

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG012918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | ---Internal Standard--- |       |        |      |
|----------------------|-------|---------|-------|----------|-------------------------|-------|--------|------|
|                      |       |         |       |          | #                       | RT    | Resp   | Conc |
| Butane, 2-methoxy... | 3.46  | 20.5    | ng    | 99480    | 1                       | 8.70  | 97198  | 20.0 |
| unknown8.22          | 8.22  | 103.0   | ng    | 500445   | 1                       | 8.70  | 97198  | 20.0 |
| N-Ethyl-N-methyl-... | 15.49 | 2.7     | ng    | 39081    | 3                       | 15.34 | 293821 | 20.0 |
| 2-Ethyl-2-(p-toly... | 15.82 | 3.7     | ng    | 53852    | 3                       | 15.34 | 293821 | 20.0 |
| unknown16.16         | 16.16 | 2.8     | ng    | 40405    | 3                       | 15.34 | 293821 | 20.0 |
| 1,4-Benzenediamin... | 16.21 | 6.2     | ng    | 91511    | 3                       | 15.34 | 293821 | 20.0 |
| unknown16.69         | 16.69 | 18.9    | ng    | 277273   | 3                       | 15.34 | 293821 | 20.0 |
| unknown18.89         | 18.89 | 2.6     | ng    | 50592    | 4                       | 18.07 | 384315 | 20.0 |