

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

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-BG011218.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         | 3.464       | 25            | 30          | 49           | rVB      | 326029         | 496154        | 12.78%          | 2.782%        |
| 2         | 5.661       | 395           | 404         | 415          | rBV      | 832630         | 1474993       | 38.00%          | 8.269%        |
| 3         | 6.155       | 481           | 488         | 500          | rBV      | 437572         | 731749        | 18.85%          | 4.102%        |
| 4         | 7.835       | 766           | 774         | 787          | rBV      | 393116         | 712444        | 18.35%          | 3.994%        |
| 5         | 8.235       | 835           | 842         | 855          | rVB      | 520865         | 958670        | 24.70%          | 5.374%        |
| 6         | 8.722       | 916           | 925         | 933          | rBV2     | 101585         | 186979        | 4.82%           | 1.048%        |
| 7         | 9.057       | 972           | 982         | 989          | rBV      | 477110         | 904180        | 23.29%          | 5.069%        |
| 8         | 9.921       | 1122          | 1129        | 1140         | rVB      | 340716         | 657436        | 16.94%          | 3.686%        |
| 9         | 10.979      | 1302          | 1309        | 1316         | rBV2     | 46425          | 87234         | 2.25%           | 0.489%        |
| 10        | 11.225      | 1342          | 1351        | 1358         | rBV      | 41050          | 72038         | 1.86%           | 0.404%        |
| 11        | 11.601      | 1400          | 1415        | 1420         | rBV      | 141883         | 280257        | 7.22%           | 1.571%        |
| 12        | 11.654      | 1420          | 1424        | 1431         | rVB      | 72006          | 127642        | 3.29%           | 0.716%        |
| 13        | 12.759      | 1607          | 1612        | 1620         | rVB2     | 28560          | 49478         | 1.27%           | 0.277%        |
| 14        | 13.217      | 1683          | 1690        | 1699         | rVB      | 87254          | 151805        | 3.91%           | 0.851%        |
| 15        | 13.435      | 1717          | 1727        | 1735         | rBV      | 68552          | 127331        | 3.28%           | 0.714%        |
| 16        | 13.987      | 1813          | 1821        | 1833         | rBV      | 1352272        | 2154830       | 55.51%          | 12.080%       |
| 17        | 14.533      | 1904          | 1914        | 1923         | rBV4     | 16855          | 42053         | 1.08%           | 0.236%        |
| 18        | 14.680      | 1930          | 1939        | 1944         | rBV3     | 30841          | 62142         | 1.60%           | 0.348%        |
| 19        | 15.362      | 2049          | 2055        | 2062         | rVB2     | 264741         | 444934        | 11.46%          | 2.494%        |
| 20        | 16.831      | 2298          | 2305        | 2317         | rBV      | 1064368        | 1771519       | 45.64%          | 9.931%        |
| 21        | 18.094      | 2513          | 2520        | 2526         | rBV      | 420862         | 663261        | 17.09%          | 3.718%        |
| 22        | 20.632      | 2945          | 2952        | 2961         | rBV      | 2817794        | 3881806       | 100.00%         | 21.762%       |
| 23        | 22.424      | 3249          | 3257        | 3266         | rBV2     | 478960         | 878690        | 22.64%          | 4.926%        |
| 24        | 26.114      | 3871          | 3885        | 3899         | rBV2     | 288178         | 919888        | 23.70%          | 5.157%        |

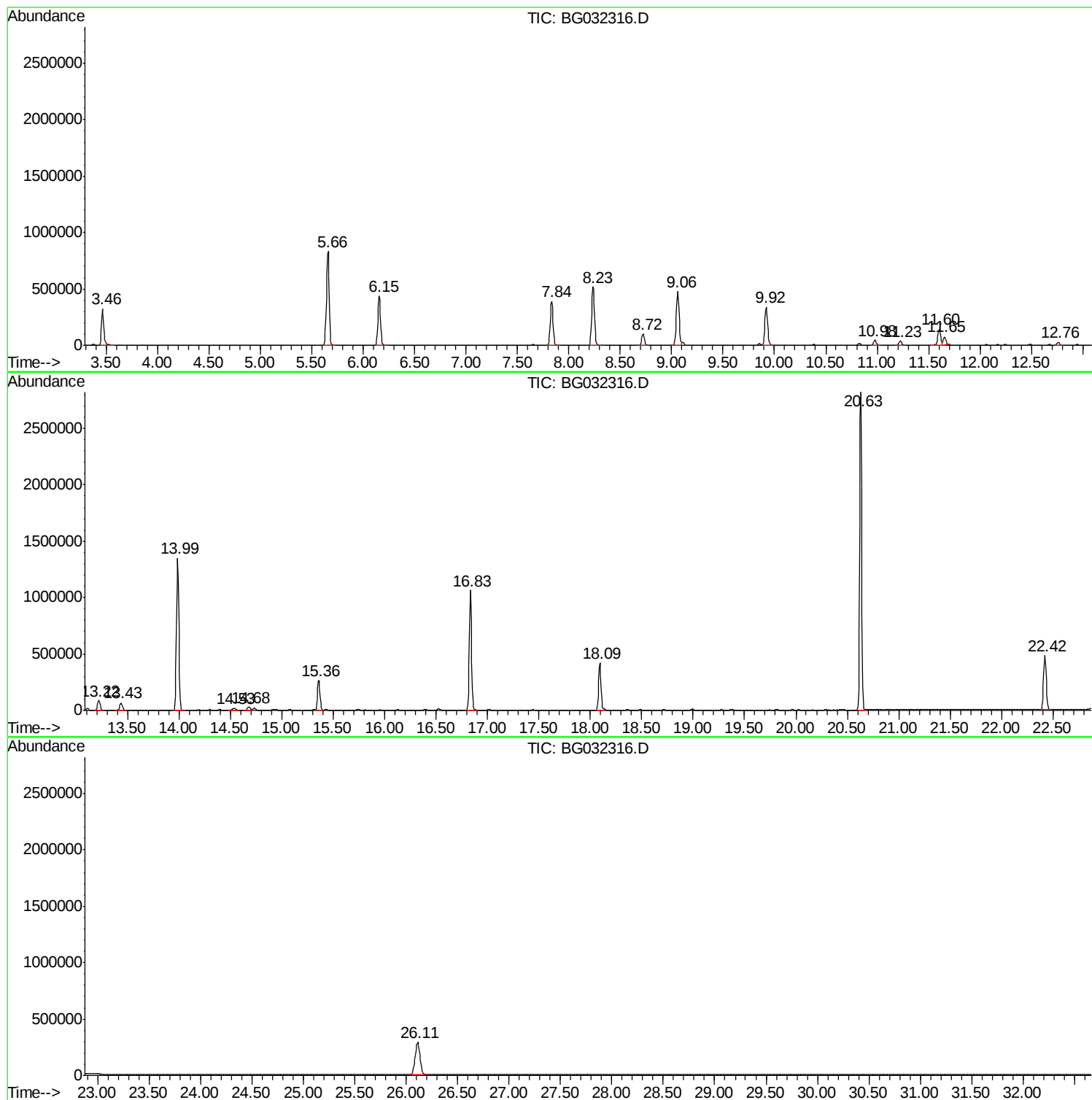
Sum of corrected areas: 17837513

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

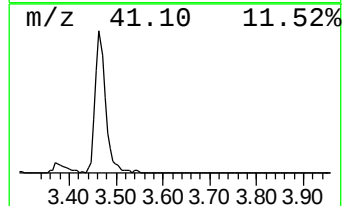
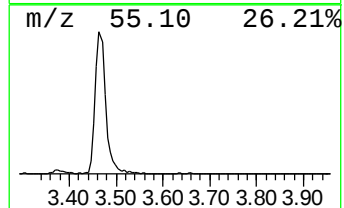
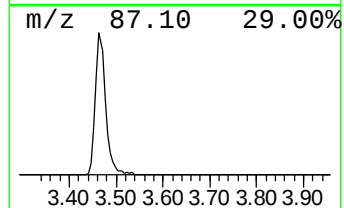
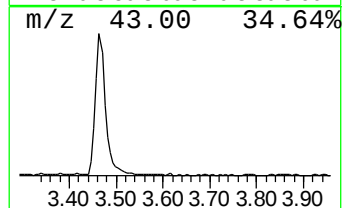
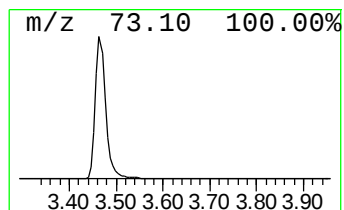
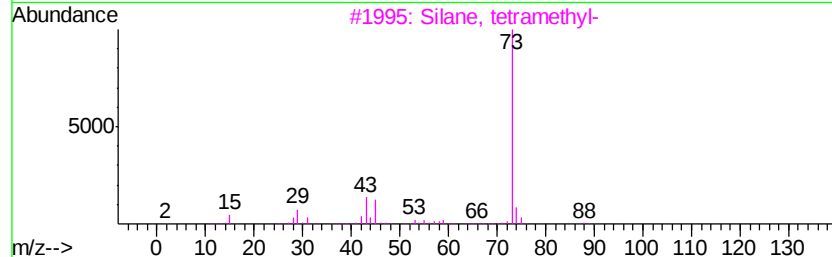
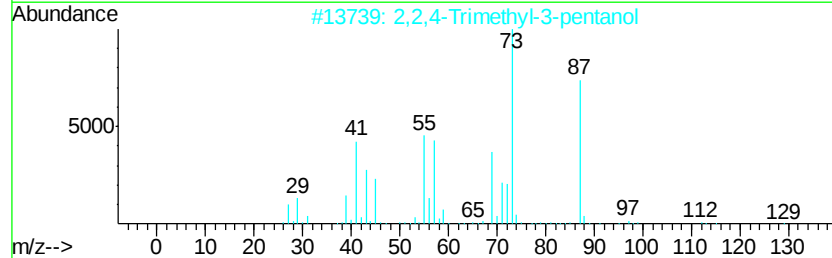
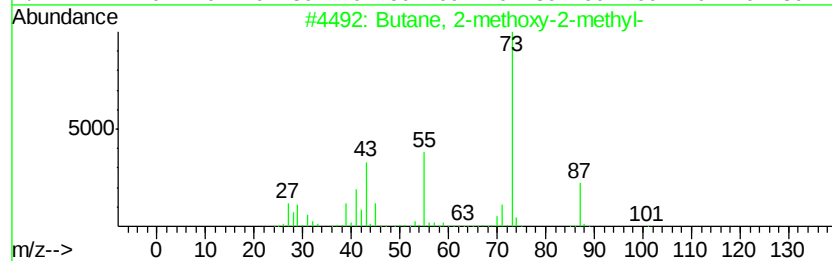
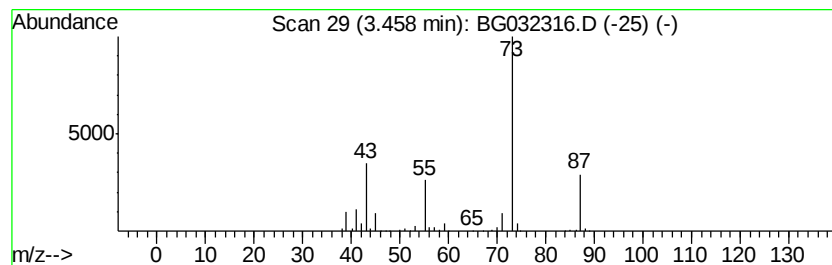
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 3.46 | 53.07 ng | 496154 | 1,4-Dichlorobenzene-d4 | 8.73 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 9    |
| 4    |    | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 5    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 9    |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

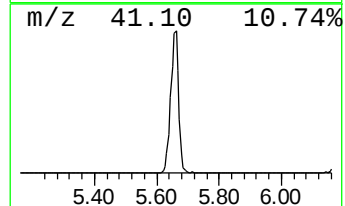
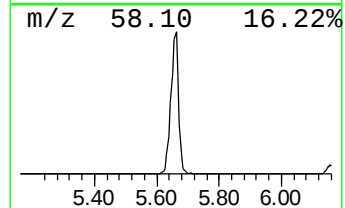
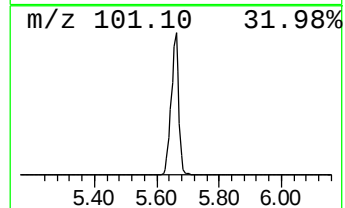
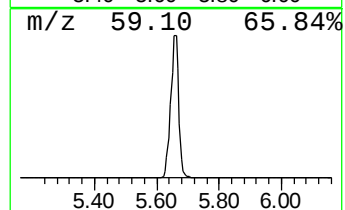
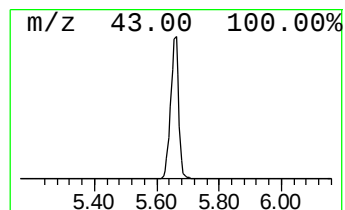
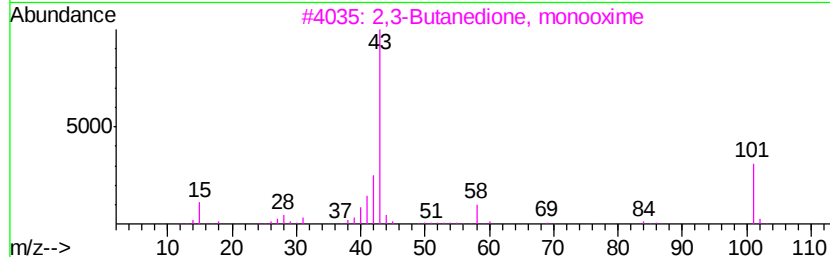
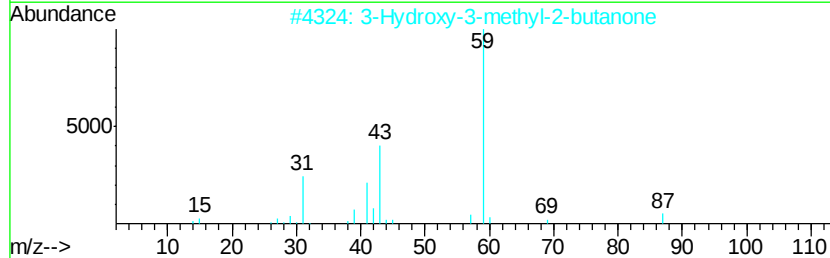
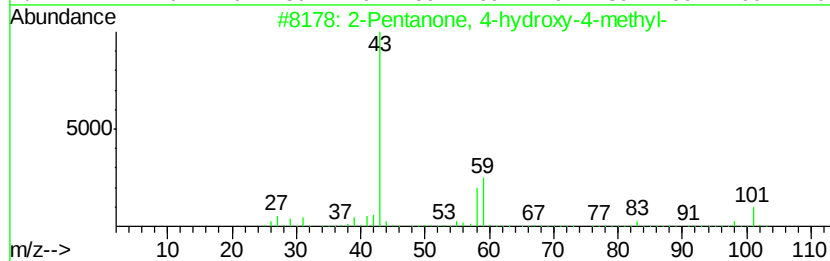
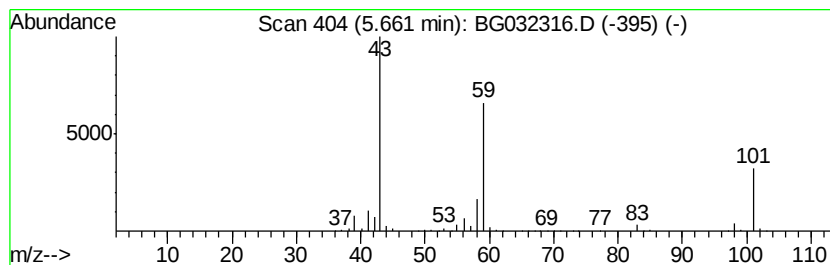
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 5.66 | 157.77 ng | 1474990 | 1,4-Dichlorobenzene-d4 | 8.73 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 64   |
| 2    |    | 3-Hydroxy-3-methyl-2-butanone    | 102 | C5H10O2 | 000115-22-0 | 25   |
| 3    |    | 2,3-Butanedione, monooxime       | 101 | C4H7NO2 | 000057-71-6 | 25   |
| 4    |    | Acetone                          | 58  | C3H6O   | 000067-64-1 | 9    |
| 5    |    | 5-Hexen-2-one                    | 98  | C6H10O  | 000109-49-9 | 9    |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

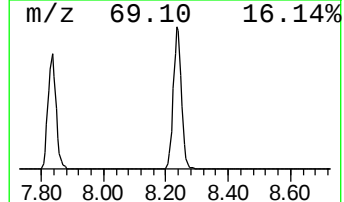
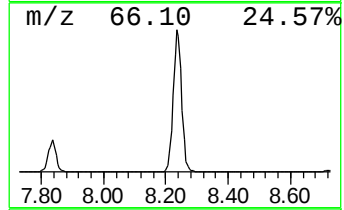
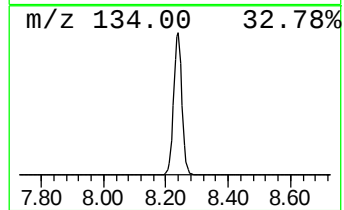
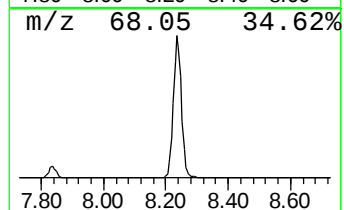
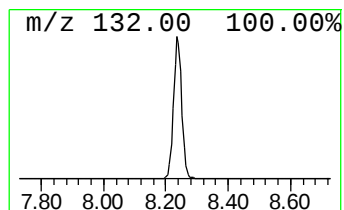
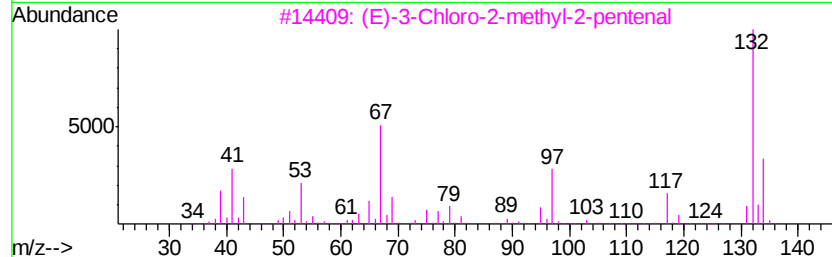
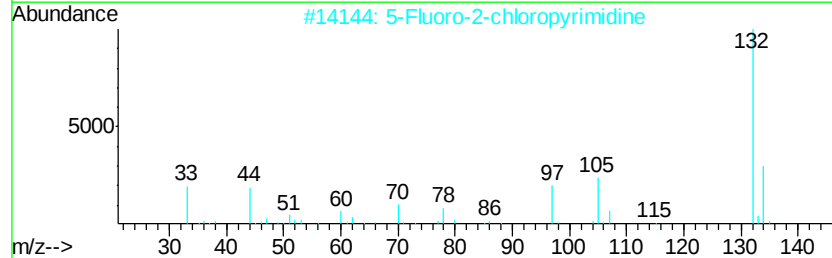
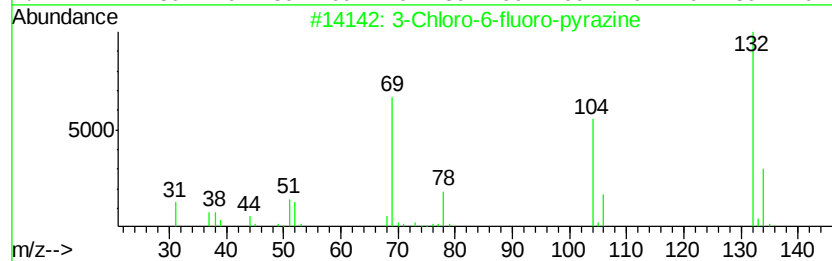
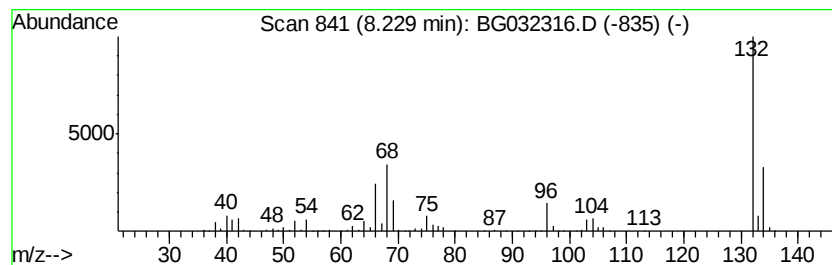
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown8.23 Concentration Rank 2

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 8.23 | 102.54 ng | 958670 | 1,4-Dichlorobenzene-d4 | 8.73 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 25   |
| 3    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 4    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 5    |    | 1,3,5-Trifluorobenzene           | 132 | C6H3F3    | 000372-38-3  | 10   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

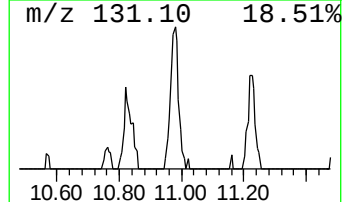
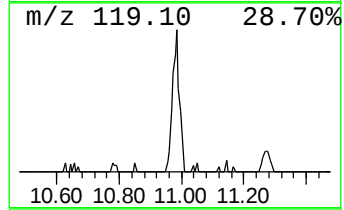
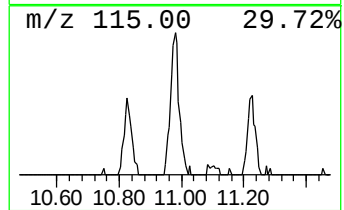
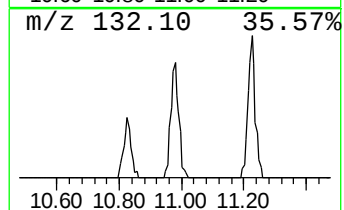
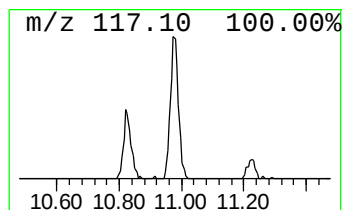
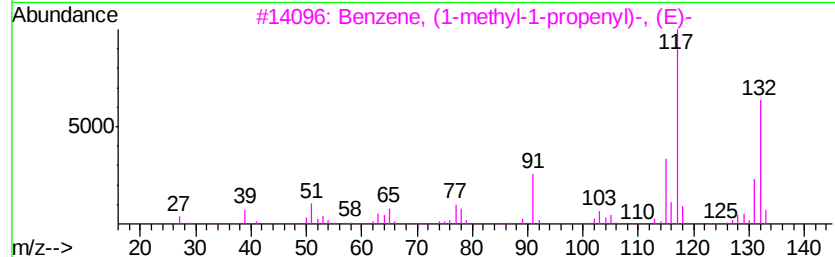
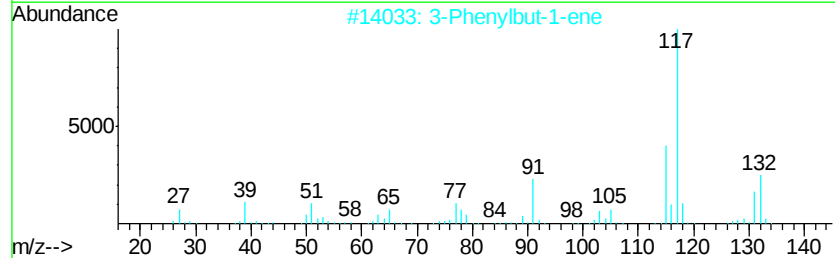
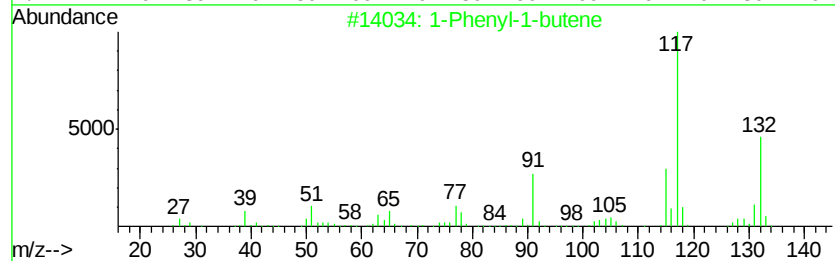
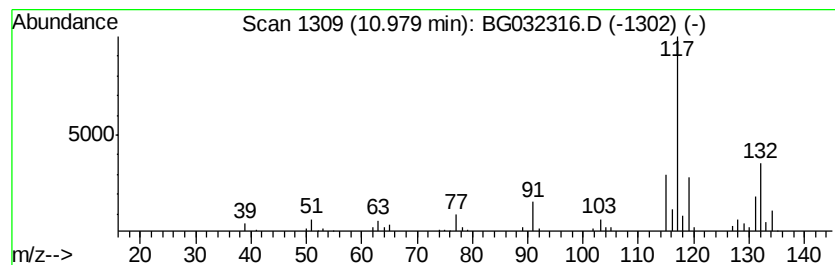
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 1-Phenyl-1-butene Concentration Rank 6

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 10.98 | 6.23 ng | 87234 | Naphthalene-d8   | 11.60 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1-Phenyl-1-butene                   | 132 | C10H12  | 000824-90-8 | 93   |
| 2    |    | 3-Phenylbut-1-ene                   | 132 | C10H12  | 000934-10-1 | 87   |
| 3    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 87   |
| 4    |    | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 81   |
| 5    |    | 1H-Indene, 2,3-dihydro-2-methyl-    | 132 | C10H12  | 000824-63-5 | 80   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

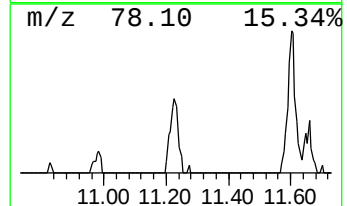
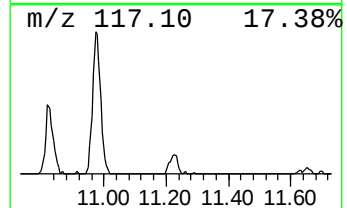
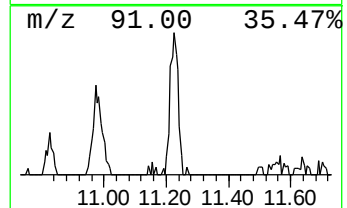
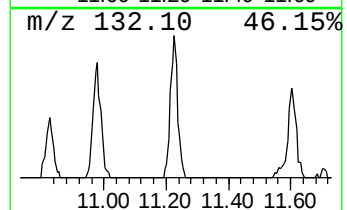
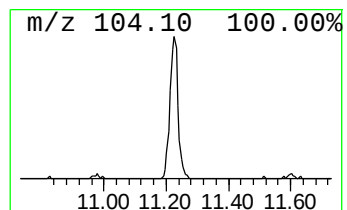
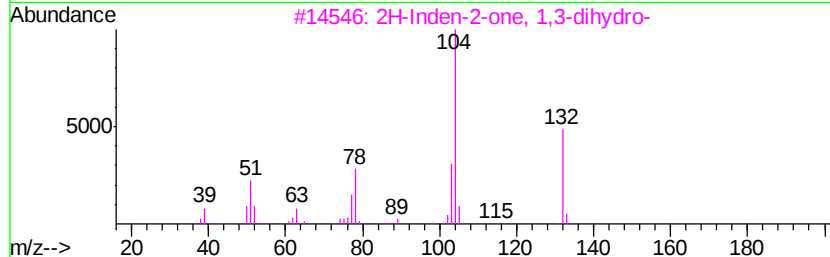
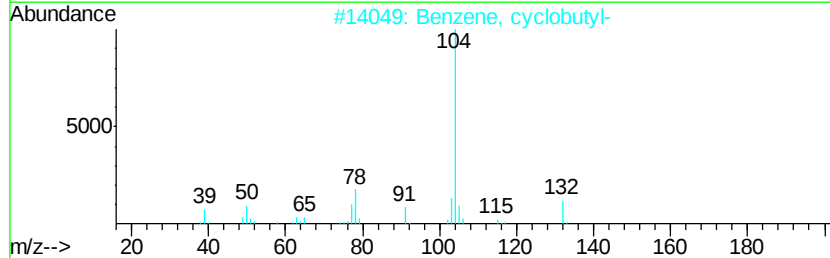
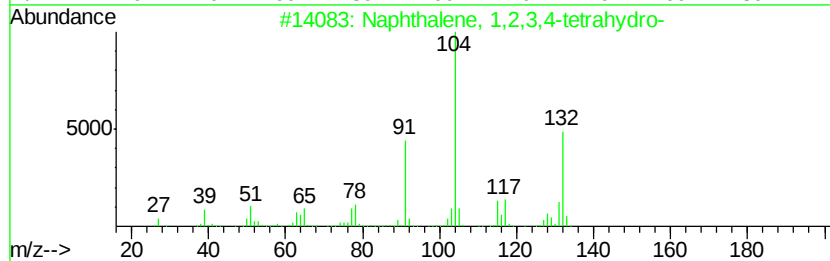
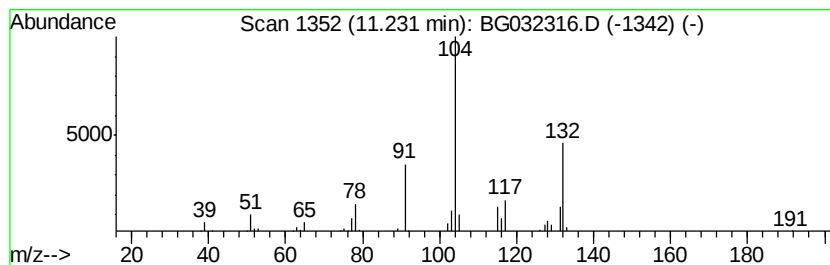
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 7

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.23 | 5.14 ng | 72038 | Naphthalene-d8   | 11.60 |

| Hit# of 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|-----------|----------------------------------|-----|---------|-------------|------|
| 1         | Naphthalene, 1,2,3,4-tetrahydro- | 132 | C10H12  | 000119-64-2 | 93   |
| 2         | Benzene, cyclobutyl-             | 132 | C10H12  | 004392-30-7 | 53   |
| 3         | 2H-Inden-2-one, 1,3-dihydro-     | 132 | C9H8O   | 000615-13-4 | 52   |
| 4         | Indan, epoxide                   | 132 | C9H8O   | 000768-22-9 | 50   |
| 5         | 2-Methylbenzyl cyanide           | 131 | C9H9N   | 022364-68-7 | 46   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
IGW-5

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

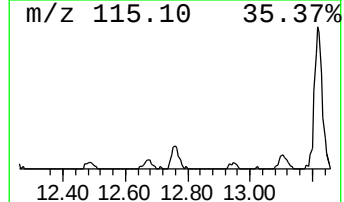
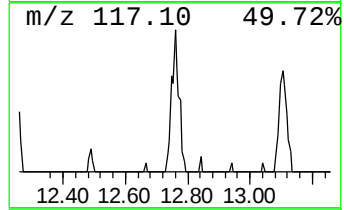
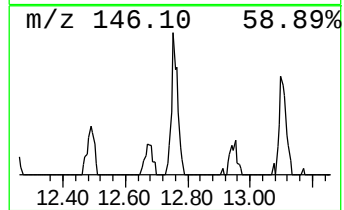
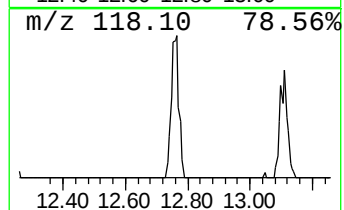
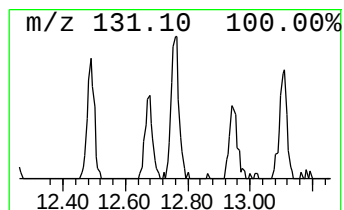
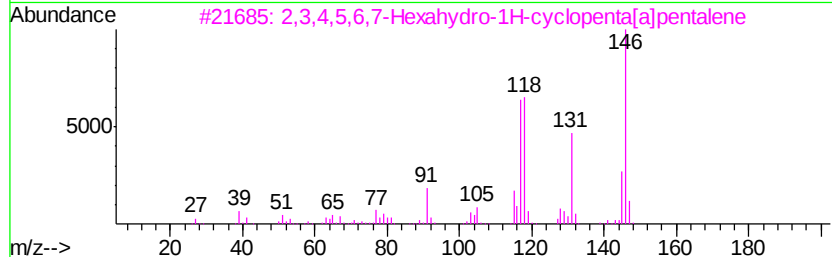
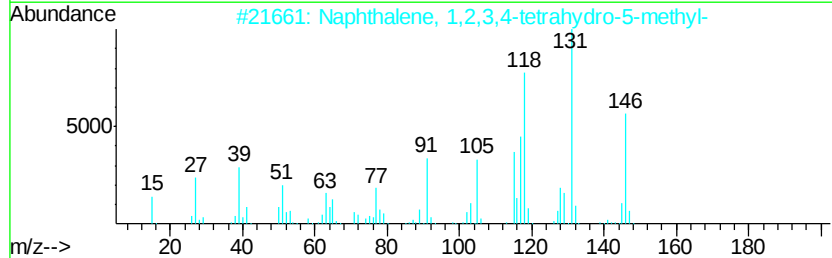
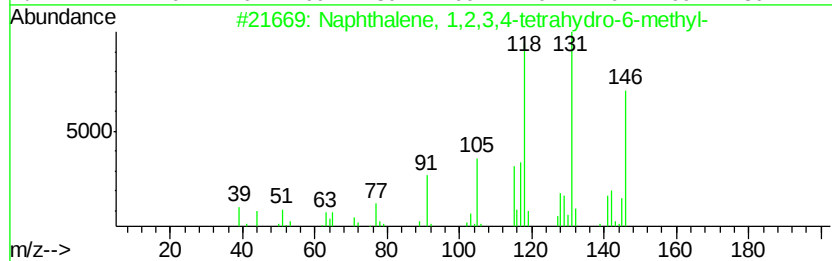
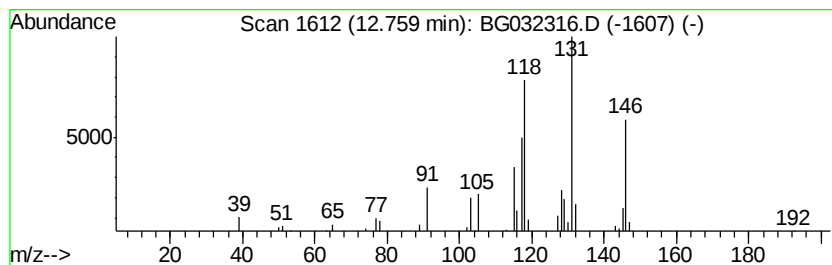
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 12.76 | 3.53 ng | 49478 | Naphthalene-d8   | 11.60 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9  | 95   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5  | 94   |
| 3    |    | 2,3,4,5,6,7-Hexahydro-1H-cyclope... | 146 | C11H14  | 1000189-31-0 | 70   |
| 4    |    | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2  | 53   |
| 5    |    | Benzene, (2-methyl-1-methylenepr... | 146 | C11H14  | 017498-71-4  | 50   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

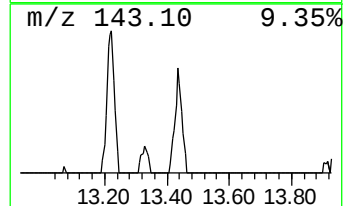
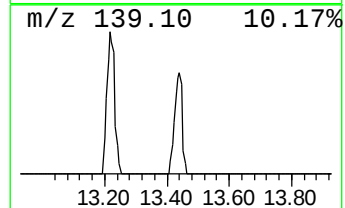
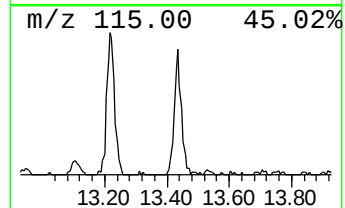
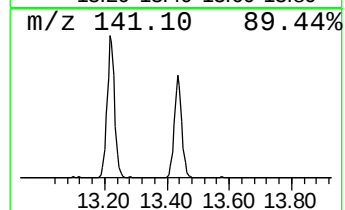
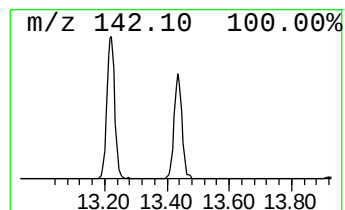
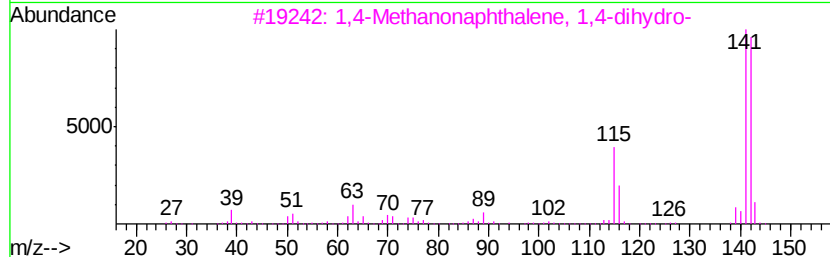
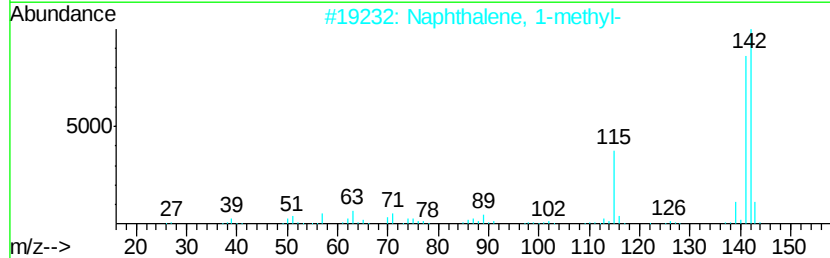
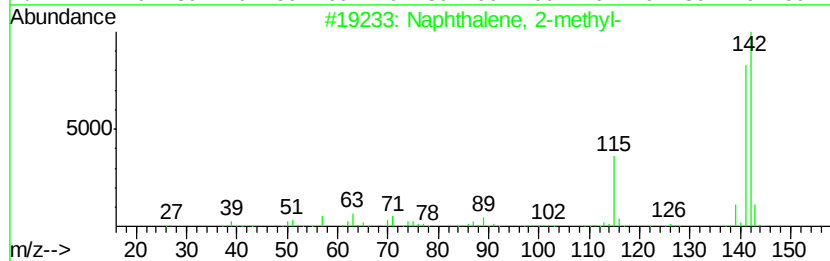
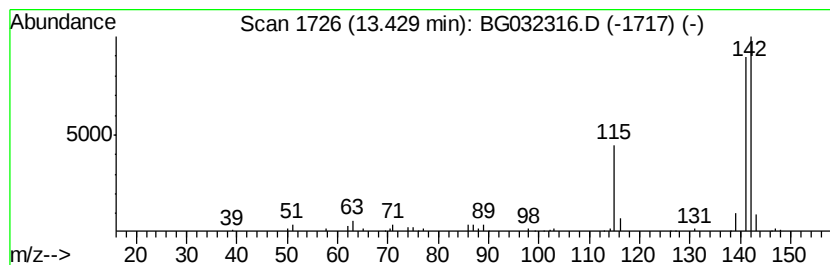
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.43 | 9.09 ng | 127331 | Naphthalene-d8   | 11.60 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 94   |
| 2    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 93   |
| 3    |    | 1,4-Methanonaphthalene, 1,4-dihv... | 142 | C11H10  | 004453-90-1 | 91   |
| 4    |    | Benzocycloheptatriene               | 142 | C11H10  | 000264-09-5 | 91   |
| 5    |    | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 | C11H10  | 002443-46-1 | 78   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

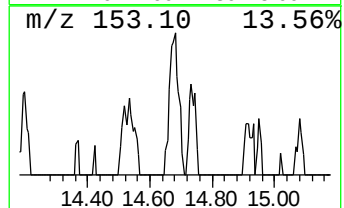
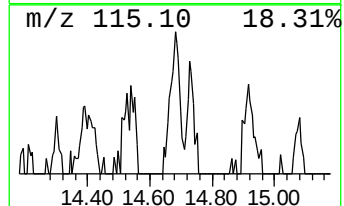
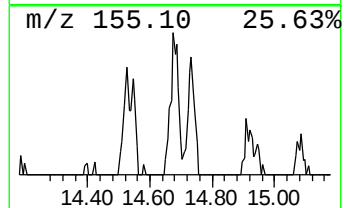
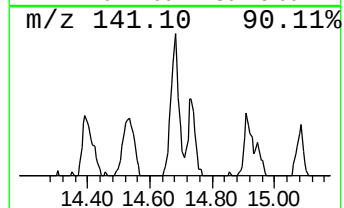
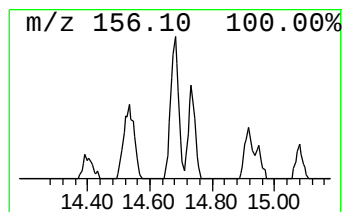
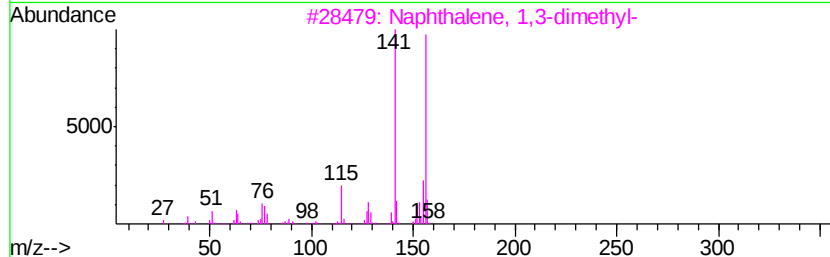
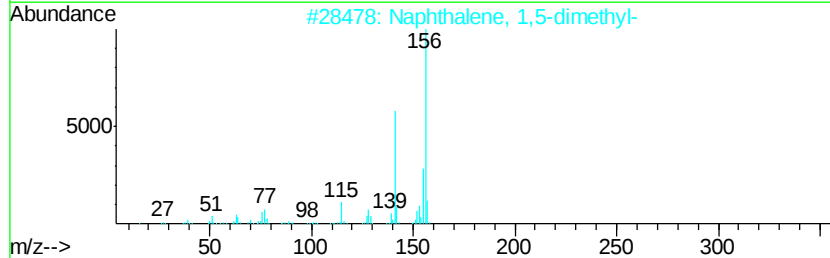
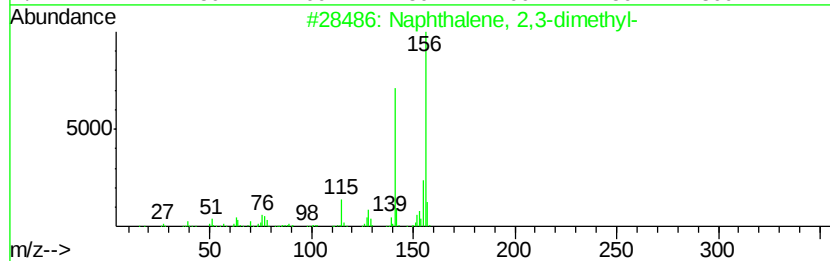
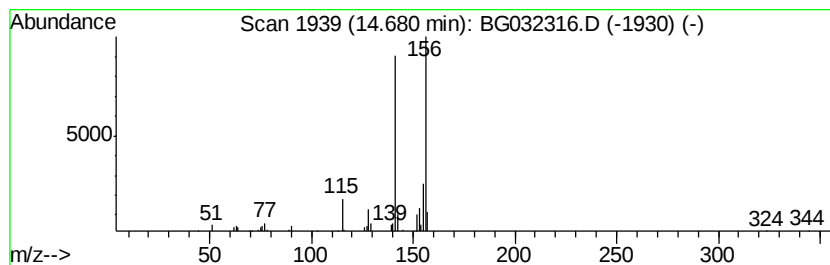
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Naphthalene, 2,3-dimethyl- Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.68 | 2.79 ng | 62142 | Acenaphthene-d10 | 15.36 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 95   |
| 2    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 93   |
| 3    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 93   |
| 4    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 93   |
| 5    |    | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 90   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG012518\  
Data File : BG032316.D  
Acq On : 26 Jan 2018 6:01  
Operator : SJ/JU  
Sample : J1267-02  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
IGW-5

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG011218.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  | 53.1    | ng    | 496154   | 1                     | 8.73  | 186979 | 20.0 |
| 2-Pentanone, 4-hy... | 5.66  | 157.8   | ng    | 1474990  | 1                     | 8.73  | 186979 | 20.0 |
| unknown8.23          | 8.23  | 102.5   | ng    | 958670   | 1                     | 8.73  | 186979 | 20.0 |
| 1-Phenyl-1-butene    | 10.98 | 6.2     | ng    | 87234    | 2                     | 11.60 | 280257 | 20.0 |
| Naphthalene, 1,2,... | 11.23 | 5.1     | ng    | 72038    | 2                     | 11.60 | 280257 | 20.0 |
| Naphthalene, 1,2,... | 12.76 | 3.5     | ng    | 49478    | 2                     | 11.60 | 280257 | 20.0 |
| Naphthalene, 1-me... | 13.43 | 9.1     | ng    | 127331   | 2                     | 11.60 | 280257 | 20.0 |
| Naphthalene, 2,3-... | 14.68 | 2.8     | ng    | 62142    | 3                     | 15.36 | 444934 | 20.0 |