

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
 Data File : BG036040.D  
 Acq On : 1 Aug 2018 19:12  
 Operator : JU/SJ  
 Sample : J4256-03  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-A27

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:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG071218.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         | 5.133       | 308           | 314         | 321          | rBV      | 39510          | 65412         | 1.75%           | 0.370%        |
| 2         | 5.603       | 387           | 394         | 411          | rBV      | 314698         | 547191        | 14.66%          | 3.099%        |
| 3         | 7.189       | 654           | 664         | 675          | rBV      | 199388         | 391570        | 10.49%          | 2.218%        |
| 4         | 7.559       | 719           | 727         | 744          | rBV      | 581649         | 1045543       | 28.02%          | 5.922%        |
| 5         | 8.023       | 799           | 806         | 813          | rBV      | 132127         | 226370        | 6.07%           | 1.282%        |
| 6         | 8.346       | 853           | 861         | 866          | rBV      | 552469         | 984414        | 26.38%          | 5.575%        |
| 7         | 8.393       | 866           | 869         | 876          | rVB      | 111893         | 173312        | 4.64%           | 0.982%        |
| 8         | 8.511       | 883           | 889         | 895          | rBV2     | 49857          | 86002         | 2.30%           | 0.487%        |
| 9         | 8.663       | 908           | 915         | 918          | rBV4     | 28763          | 45287         | 1.21%           | 0.256%        |
| 10        | 8.710       | 918           | 923         | 929          | rVB2     | 35967          | 62223         | 1.67%           | 0.352%        |
| 11        | 9.127       | 984           | 994         | 998          | rBV      | 139341         | 247326        | 6.63%           | 1.401%        |
| 12        | 9.186       | 998           | 1004        | 1018         | rVB2     | 437648         | 979583        | 26.25%          | 5.548%        |
| 13        | 9.644       | 1077          | 1082        | 1089         | rVB      | 70200          | 121182        | 3.25%           | 0.686%        |
| 14        | 10.015      | 1139          | 1145        | 1151         | rBV6     | 26646          | 61599         | 1.65%           | 0.349%        |
| 15        | 10.220      | 1173          | 1180        | 1189         | rBV2     | 266372         | 494679        | 13.26%          | 2.802%        |
| 16        | 10.825      | 1271          | 1283        | 1288         | rBV      | 164729         | 364756        | 9.77%           | 2.066%        |
| 17        | 10.866      | 1288          | 1290        | 1297         | rVV5     | 33145          | 58258         | 1.56%           | 0.330%        |
| 18        | 10.937      | 1297          | 1302        | 1309         | rVB3     | 34804          | 65279         | 1.75%           | 0.370%        |
| 19        | 11.283      | 1355          | 1361        | 1369         | rVB      | 48315          | 88066         | 2.36%           | 0.499%        |
| 20        | 11.401      | 1375          | 1381        | 1386         | rBV      | 50787          | 93071         | 2.49%           | 0.527%        |
| 21        | 11.724      | 1431          | 1436        | 1443         | rVB3     | 19758          | 37886         | 1.02%           | 0.215%        |
| 22        | 11.930      | 1465          | 1471        | 1477         | rVB4     | 23936          | 40733         | 1.09%           | 0.231%        |
| 23        | 12.012      | 1477          | 1485        | 1491         | rBV2     | 69153          | 123190        | 3.30%           | 0.698%        |
| 24        | 12.364      | 1538          | 1545        | 1551         | rVB2     | 116975         | 192540        | 5.16%           | 1.090%        |
| 25        | 12.693      | 1595          | 1601        | 1611         | rVB3     | 69325          | 148878        | 3.99%           | 0.843%        |
| 26        | 13.269      | 1692          | 1699        | 1706         | rBV      | 1305813        | 2017628       | 54.07%          | 11.427%       |
| 27        | 13.980      | 1811          | 1820        | 1824         | rBV3     | 76714          | 156457        | 4.19%           | 0.886%        |
| 28        | 14.091      | 1834          | 1839        | 1845         | rBV      | 94705          | 142340        | 3.81%           | 0.806%        |
| 29        | 14.644      | 1926          | 1933        | 1940         | rBV2     | 324925         | 563610        | 15.10%          | 3.192%        |
| 30        | 14.708      | 1940          | 1944        | 1952         | rVB      | 52507          | 85392         | 2.29%           | 0.484%        |
| 31        | 16.136      | 2180          | 2187        | 2200         | rVV      | 1122826        | 1795616       | 48.12%          | 10.170%       |
| 32        | 16.365      | 2223          | 2226        | 2235         | rVB2     | 23726          | 41150         | 1.10%           | 0.233%        |
| 33        | 17.387      | 2391          | 2400        | 2406         | rBV      | 441533         | 681642        | 18.27%          | 3.861%        |
| 34        | 18.274      | 2545          | 2551        | 2557         | rBV4     | 34721          | 68177         | 1.83%           | 0.386%        |

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

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:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG071218.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |         |
|----|--------|------|------|------|------|---------|---------|---------|---------|
| 35 | 19.995 | 2838 | 2844 | 2854 | rBV  | 2798339 | 3731701 | 100.00% | 21.135% |
| 36 | 21.652 | 3120 | 3126 | 3136 | rVB  | 500255  | 824735  | 22.10%  | 4.671%  |
| 37 | 24.847 | 3658 | 3670 | 3681 | rBV2 | 279175  | 803436  | 21.53%  | 4.550%  |

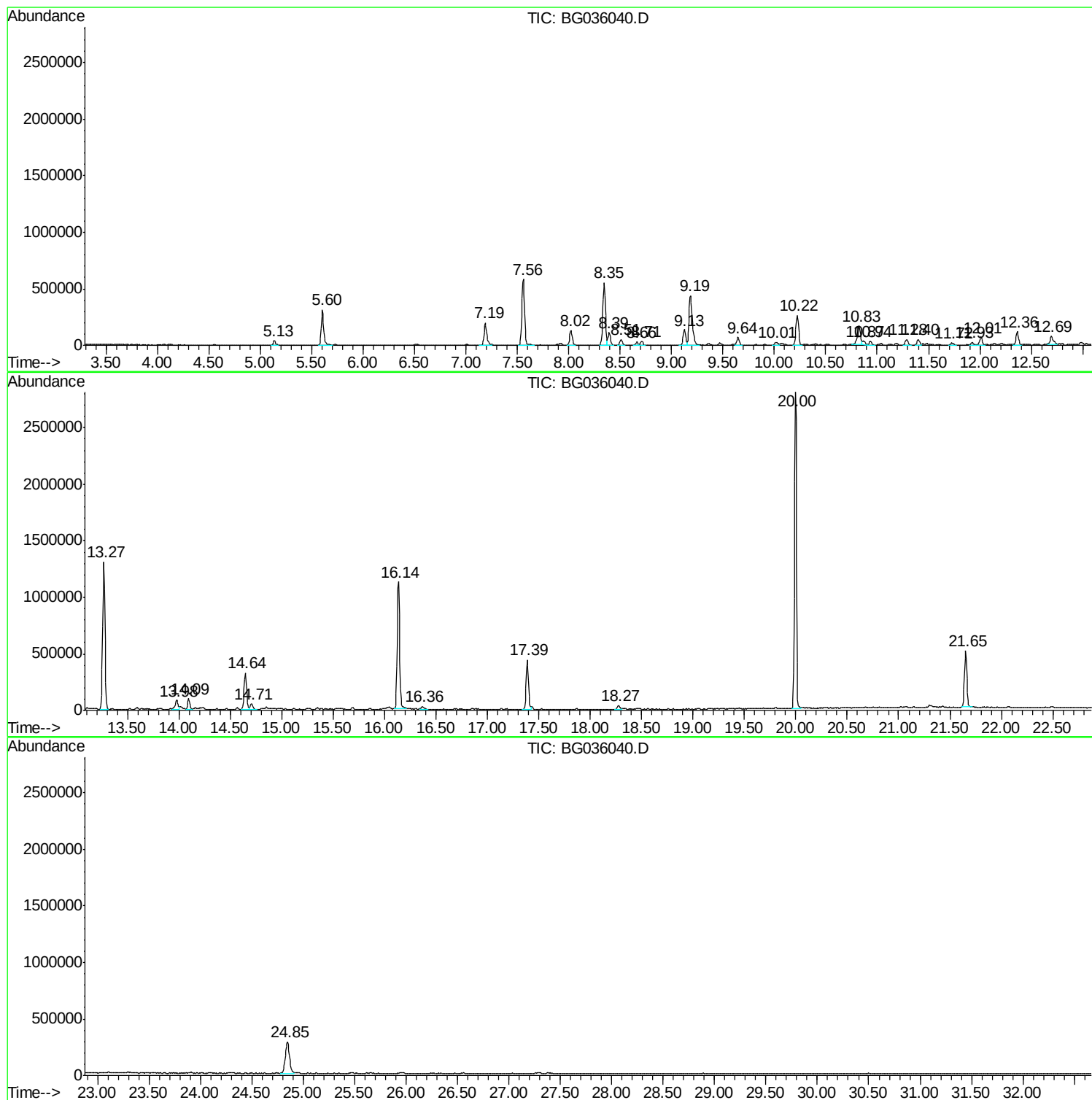
Sum of corrected areas: 17656234

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

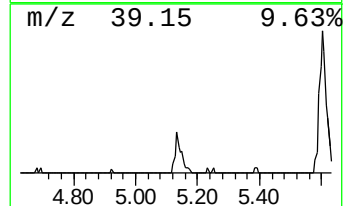
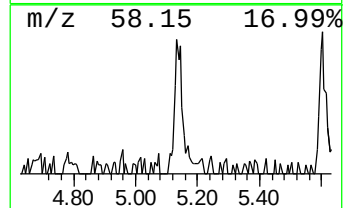
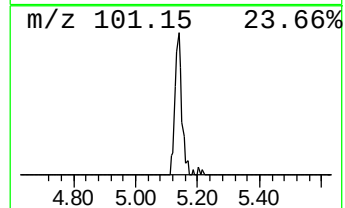
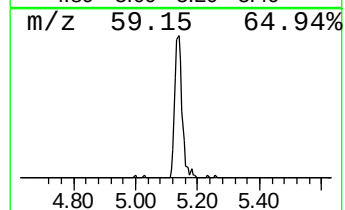
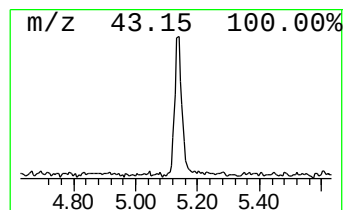
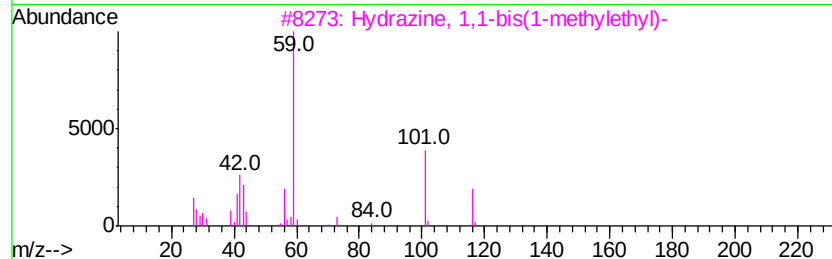
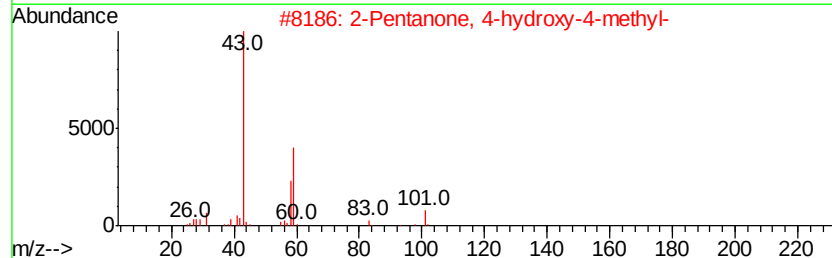
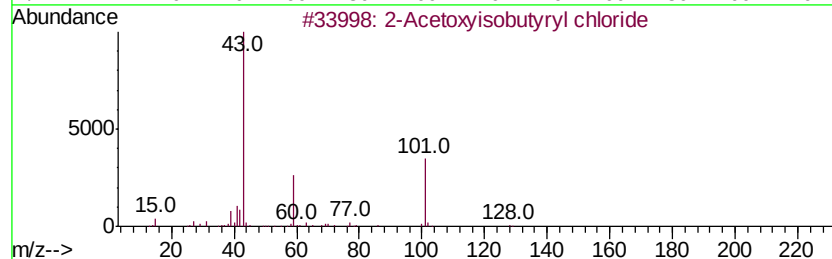
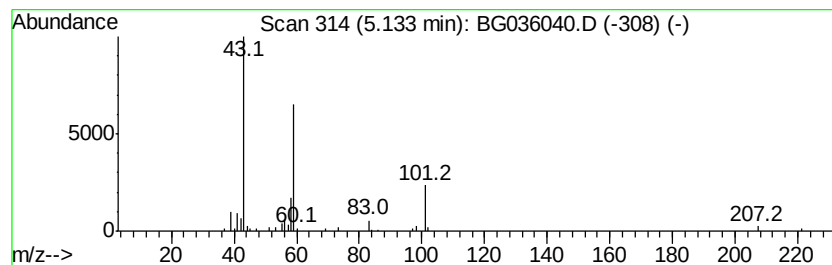
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 2-Acetoxyisobutryl chloride Concentration Rank 11

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 5.13 | 5.78 ng | 65412 | 1,4-Dichlorobenzene-d4 | 8.02 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 2-Acetoxyisobutryl chloride         | 164 | C6H9ClO3 | 040635-66-3  | 50   |
| 2    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2  | 45   |
| 3    |    |   | Hydrazine, 1,1-bis(1-methylethyl)-  | 116 | C6H16N2  | 000921-14-2  | 38   |
| 4    |    |   | n-Propyl t-butyl ether              | 116 | C7H16O   | 1000334-81-9 | 28   |
| 5    |    |   | 1,3-Dioxolane-2-methanol, 2,4-di... | 132 | C6H12O3  | 053951-43-2  | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

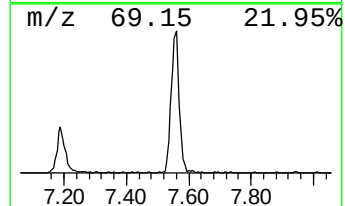
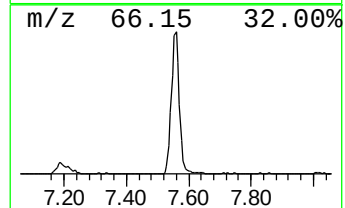
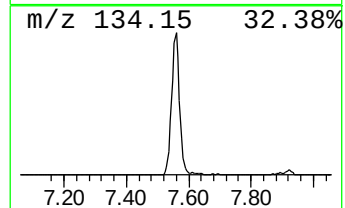
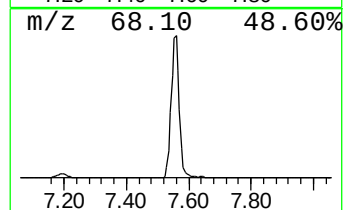
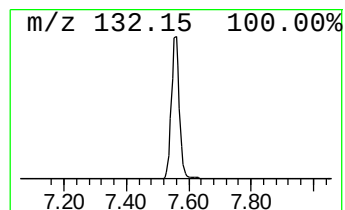
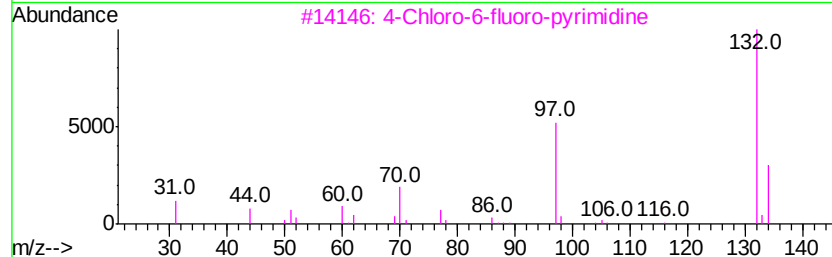
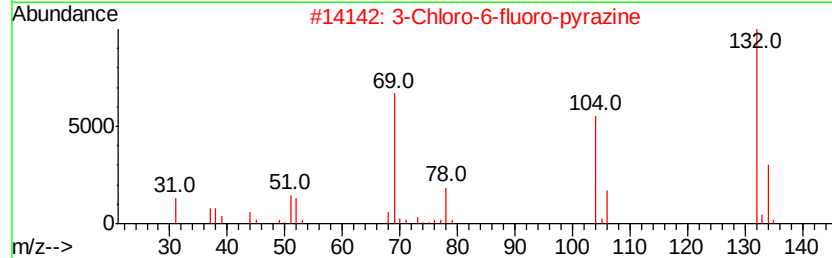
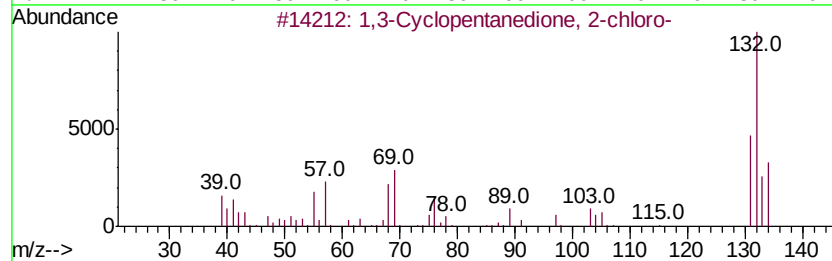
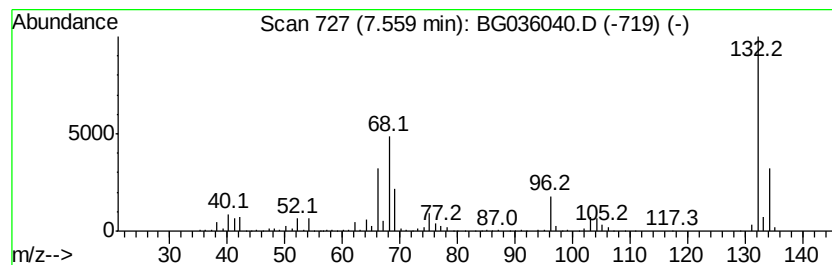
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 unknown7.56 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 7.56 | 92.37 ng | 1045540 | 1,4-Dichlorobenzene-d4 | 8.02 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 1,3-Cyclopentanedione, 2-chloro- | 132 | C5H5ClO2  | 014203-19-1  | 35   |
| 2    |    | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 3    |    | 4-Chloro-6-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 4    |    | Imidazole, 4,5-dicyano-1-methyl- | 132 | C6H4N4    | 019485-35-9  | 25   |
| 5    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

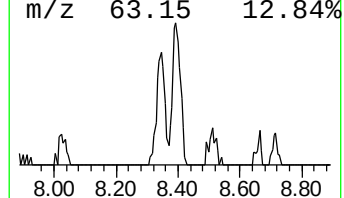
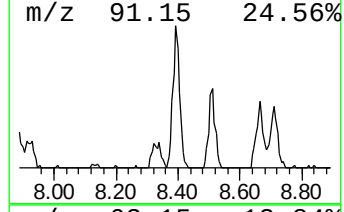
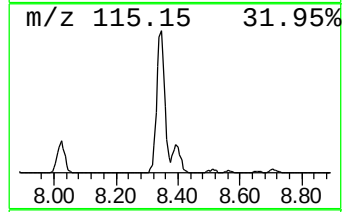
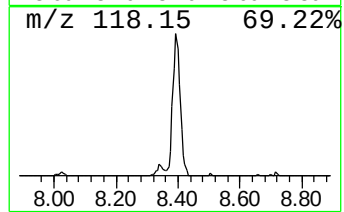
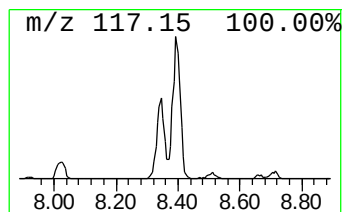
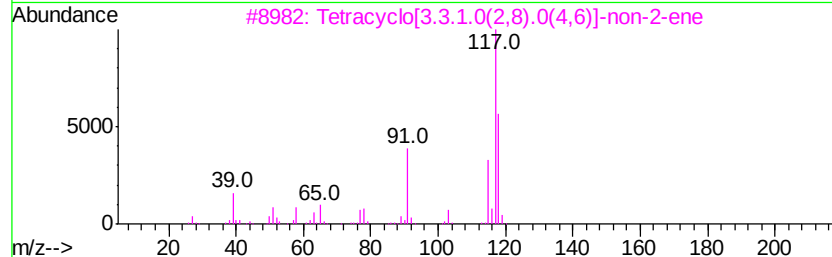
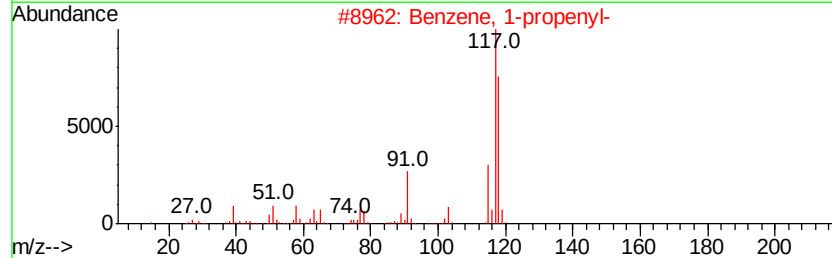
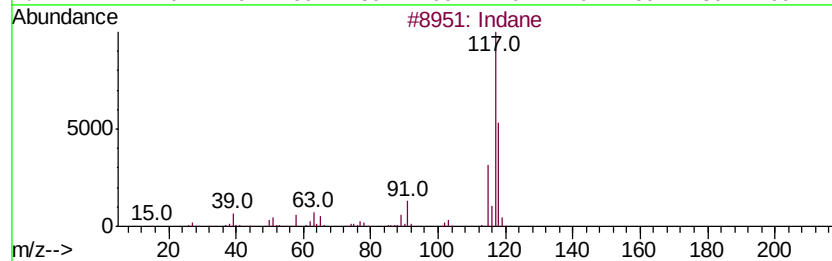
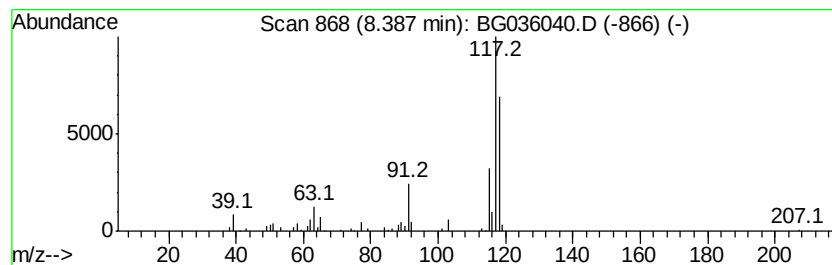
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Indane Concentration Rank 5

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 8.39 | 15.31 ng | 173312 | 1,4-Dichlorobenzene-d4 | 8.02 |

| Hit# of | 5                                | Tentative ID | MW  | MolForm | CAS#         | Qual |
|---------|----------------------------------|--------------|-----|---------|--------------|------|
| 1       | Indane                           |              | 118 | C9H10   | 000496-11-7  | 74   |
| 2       | Benzene, 1-propenyl-             |              | 118 | C9H10   | 000637-50-3  | 64   |
| 3       | Tetracyclo[3.3.1.0(2,8).0(4,6)]- | ...          | 118 | C9H10   | 1000191-13-7 | 64   |
| 4       | Benzene, cyclopropyl-            |              | 118 | C9H10   | 000873-49-4  | 64   |
| 5       | Benzene, 2-propenyl-             |              | 118 | C9H10   | 000300-57-2  | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

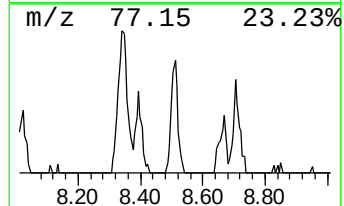
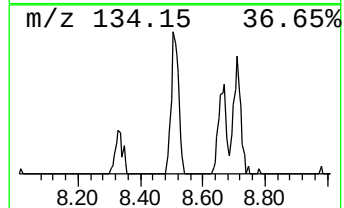
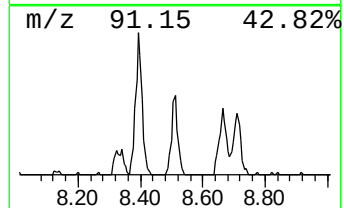
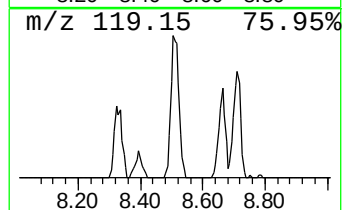
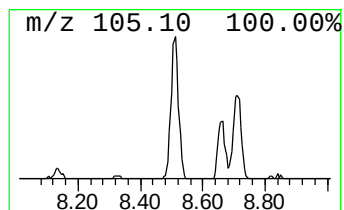
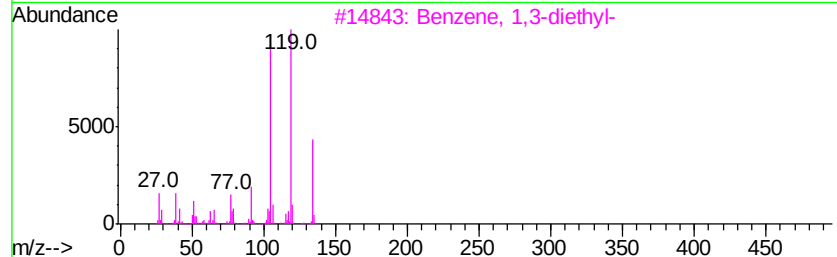
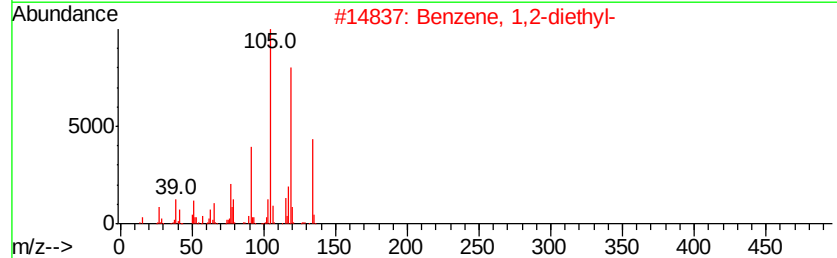
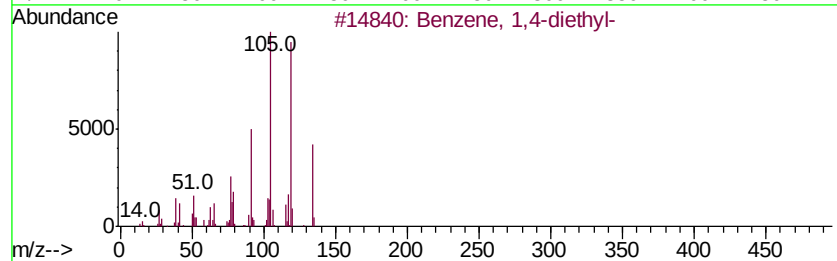
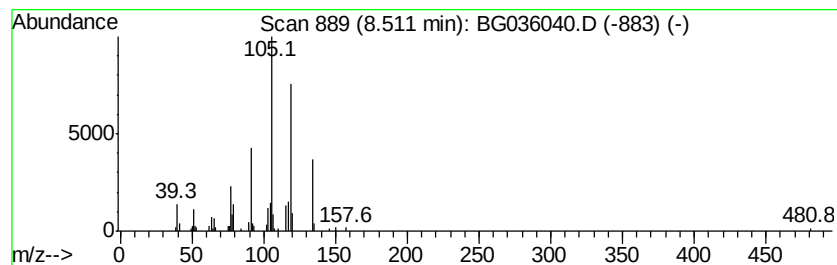
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Benzene, 1,4-diethyl- Concentration Rank 8

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 8.51 | 7.60 ng | 86002 | 1,4-Dichlorobenzene-d4 | 8.02 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,4-diethyl- | 134 | C10H14  | 000105-05-5 | 98   |
| 2    |    | Benzene, 1,2-diethyl- | 134 | C10H14  | 000135-01-3 | 96   |
| 3    |    | Benzene, 1,3-diethyl- | 134 | C10H14  | 000141-93-5 | 83   |
| 4    |    | o-Cymene              | 134 | C10H14  | 000527-84-4 | 74   |
| 5    |    | p-Cymene              | 134 | C10H14  | 000099-87-6 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

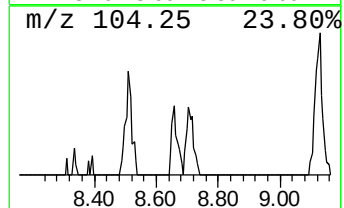
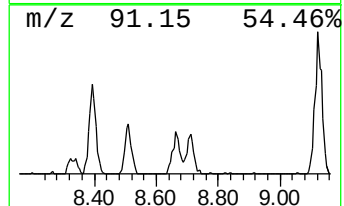
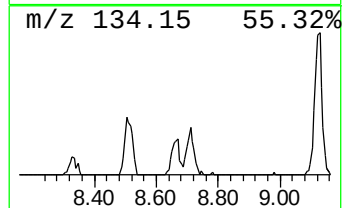
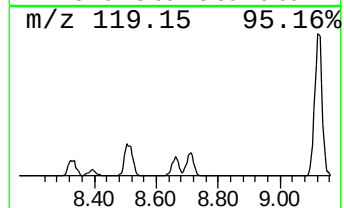
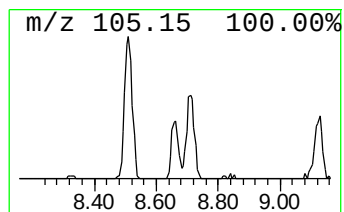
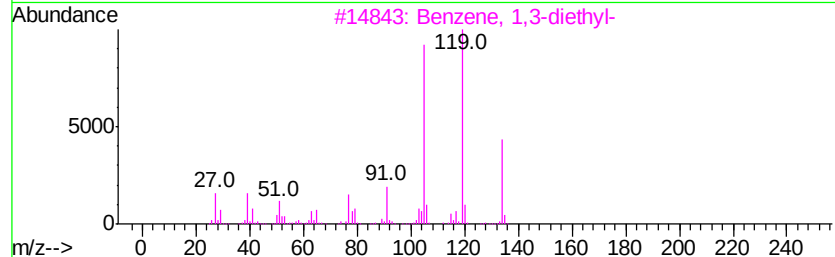
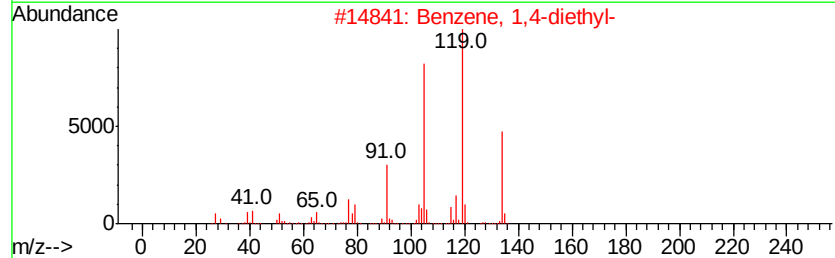
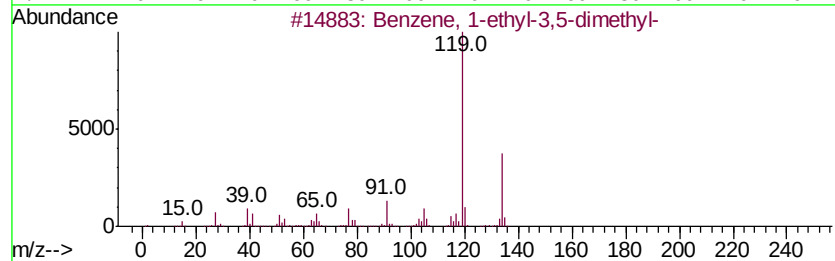
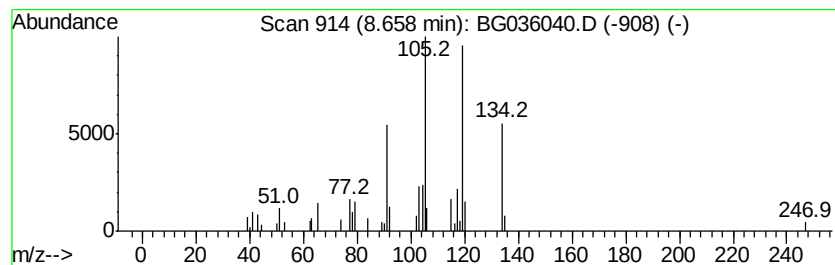
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 16

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 8.66 | 4.00 ng | 45287 | 1,4-Dichlorobenzene-d4 | 8.02 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 87   |
| 2    |    | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 87   |
| 3    |    | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 87   |
| 4    |    | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 87   |
| 5    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 87   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

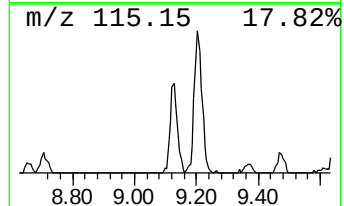
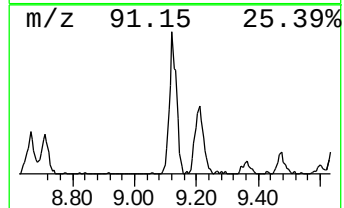
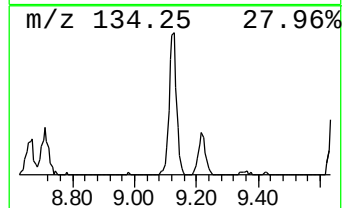
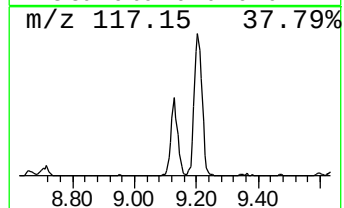
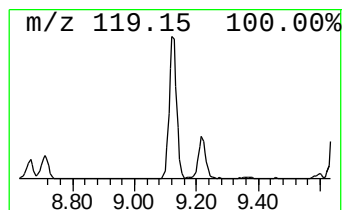
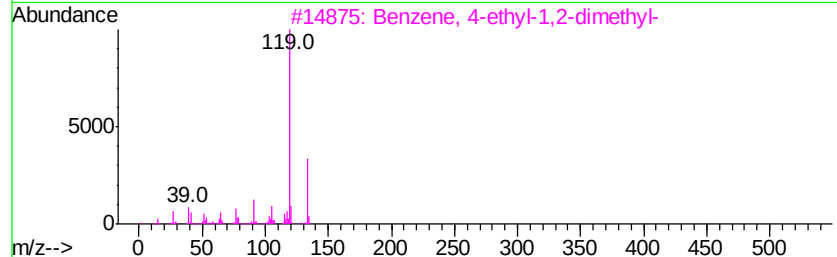
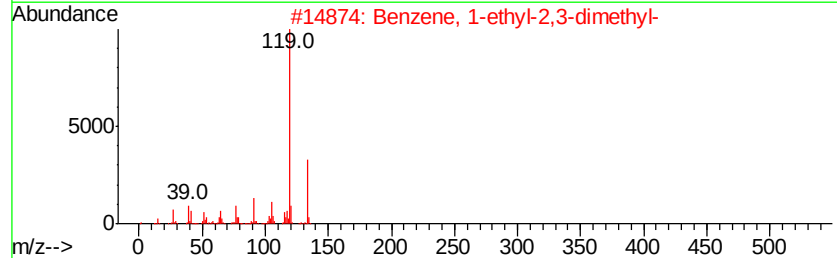
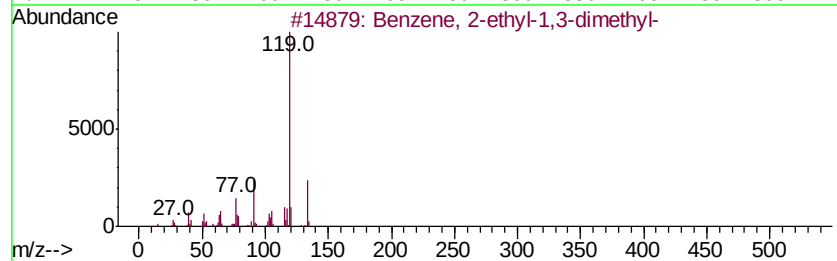
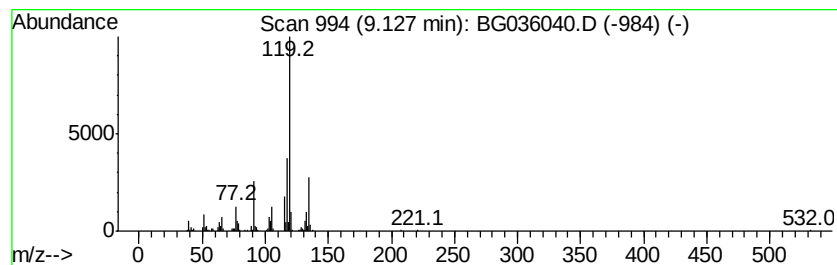
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 9.13 | 21.85 ng | 247326 | 1,4-Dichlorobenzene-d4 | 8.02 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 93   |
| 2    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 81   |
| 3    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 81   |
| 4    |    | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 76   |
| 5    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

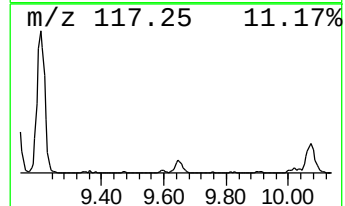
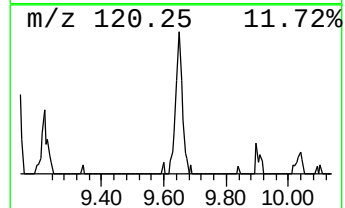
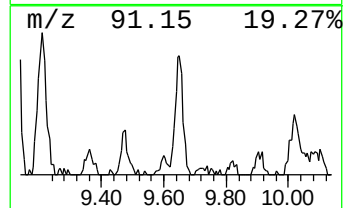
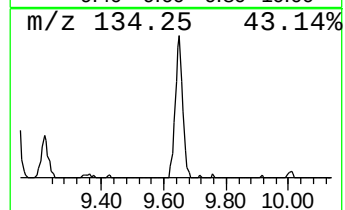
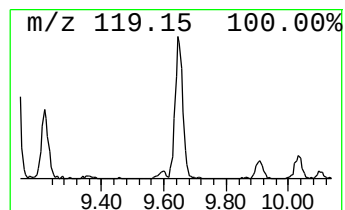
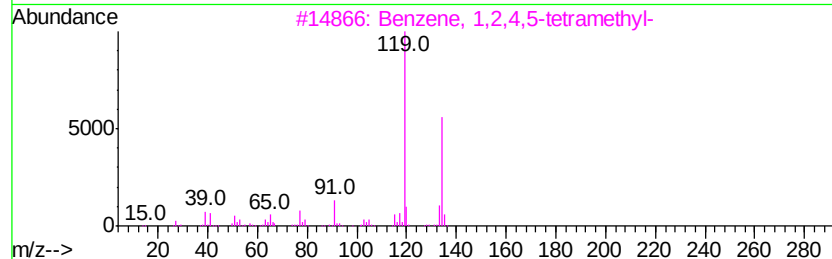
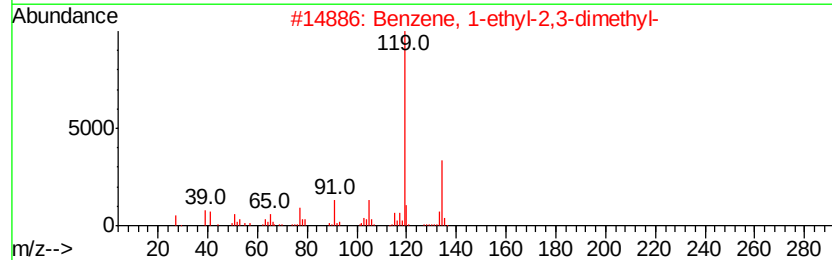
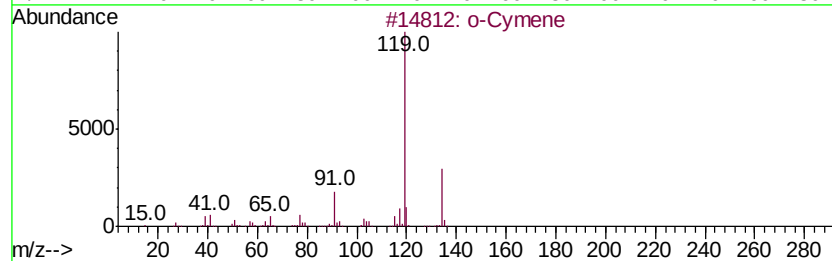
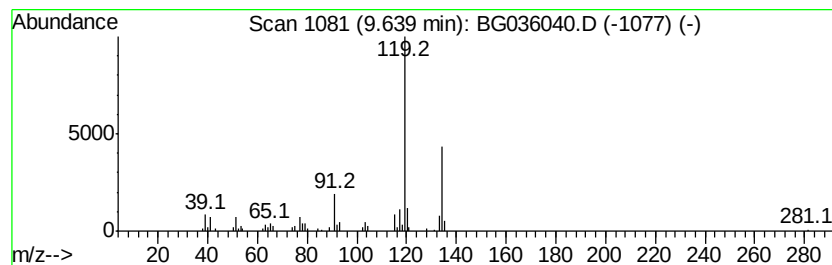
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 o-Cymene Concentration Rank 10

| R.T. | EstConc | Area   | Relative to ISTD | R.T.  |
|------|---------|--------|------------------|-------|
| 9.64 | 6.64 ng | 121182 | Naphthalene-d8   | 10.83 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 91   |
| 2    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 91   |
| 3    |    |   | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 91   |
| 4    |    |   | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 91   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

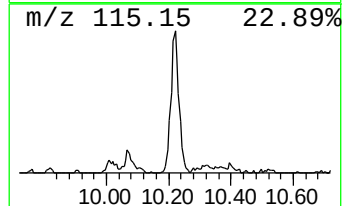
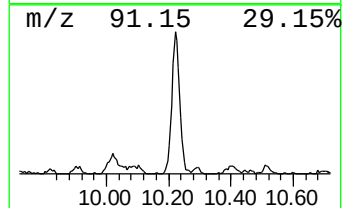
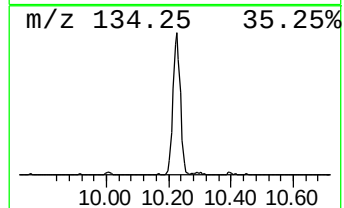
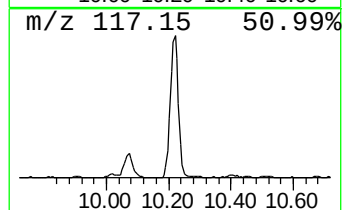
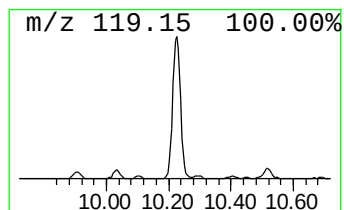
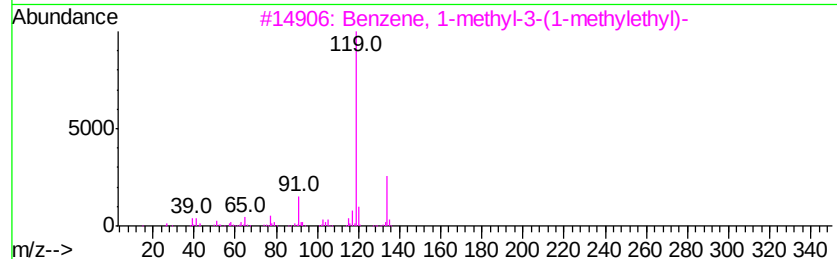
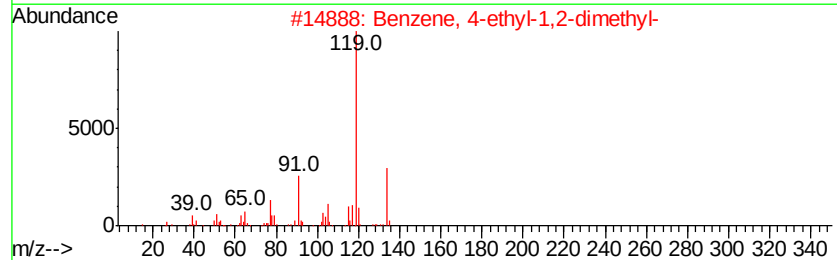
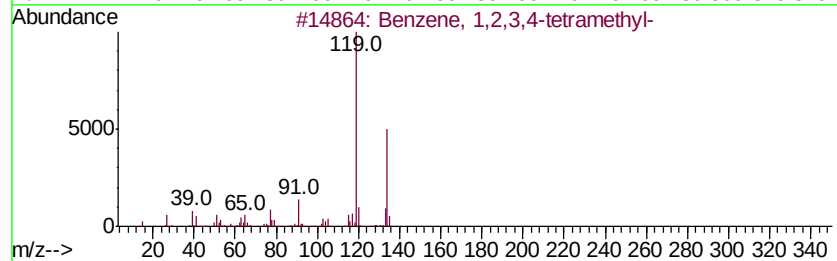
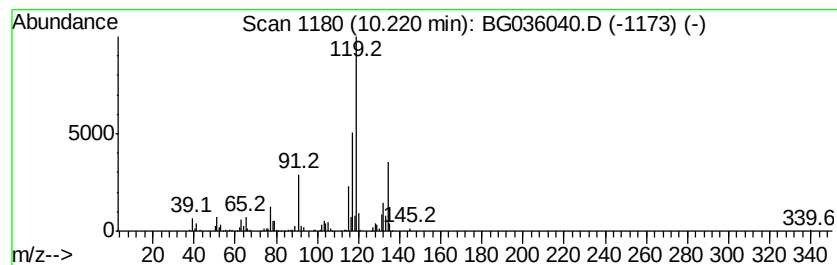
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.22 | 27.12 ng | 494679 | Napthalene-d8    | 10.83 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3,4-tetramethyl-        | 134 | C10H14  | 000488-23-3 | 64   |
| 2    |    | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 60   |
| 3    |    | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 60   |
| 4    |    | p-Cymene                             | 134 | C10H14  | 000099-87-6 | 60   |
| 5    |    | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

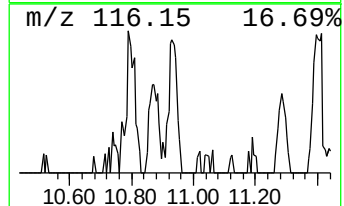
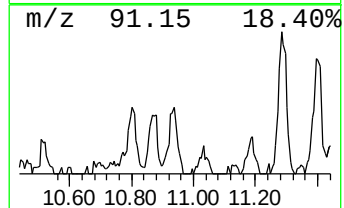
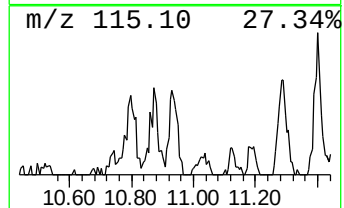
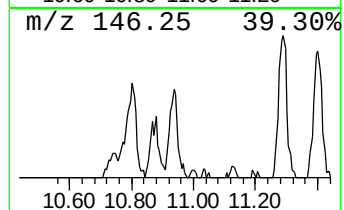
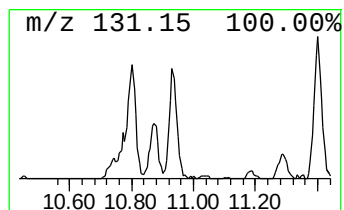
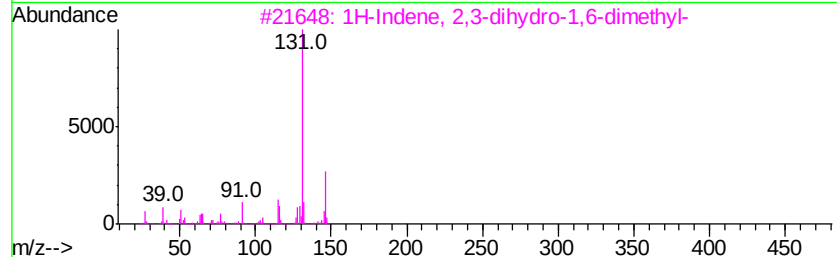
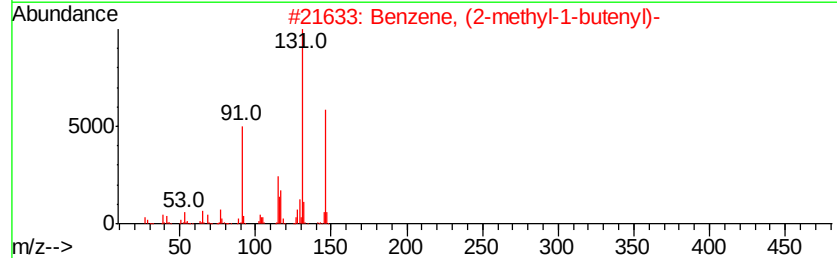
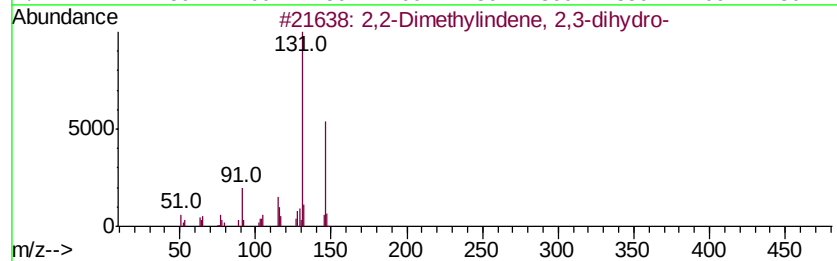
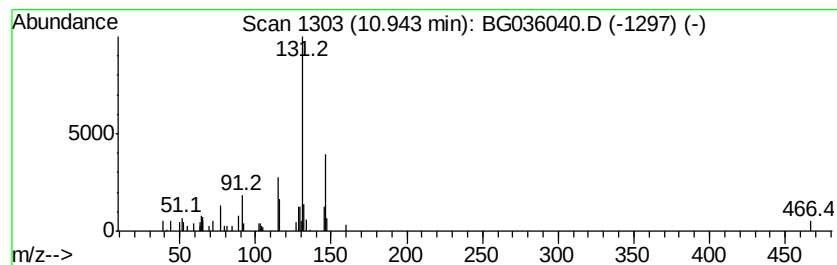
Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 2,2-Dimethylindene, 2,3-dih... Concentration Rank 17

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.        |      |
|---------|-------------------------------------|--------------|------------------|-------------|------|
| 10.94   | 3.58 ng                             | 65279        | Napthalene-d8    | 10.83       |      |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS#        | Qual |
| 1       | 2,2-Dimethylindene, 2,3-dihydro-    | 146          | C11H14           | 020836-11-7 | 91   |
| 2       | Benzene, (2-methyl-1-butenyl)-      | 146          | C11H14           | 056253-64-6 | 91   |
| 3       | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146          | C11H14           | 017059-48-2 | 90   |
| 4       | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146          | C11H14           | 006682-71-9 | 87   |
| 5       | 1H-Indene, 2,3-dihydro-5,6-dimet... | 146          | C11H14           | 001075-22-5 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

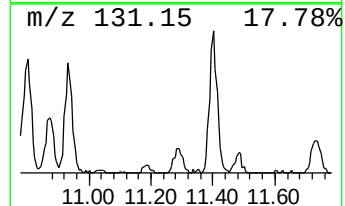
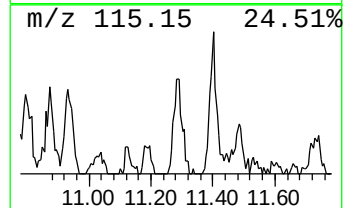
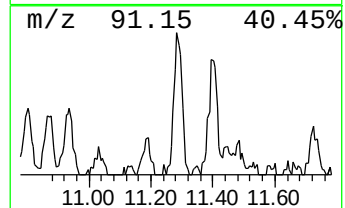
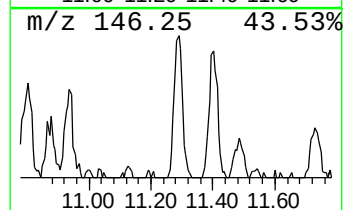
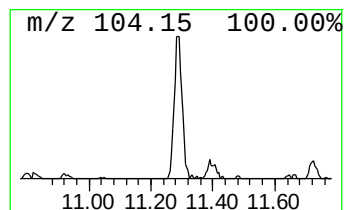
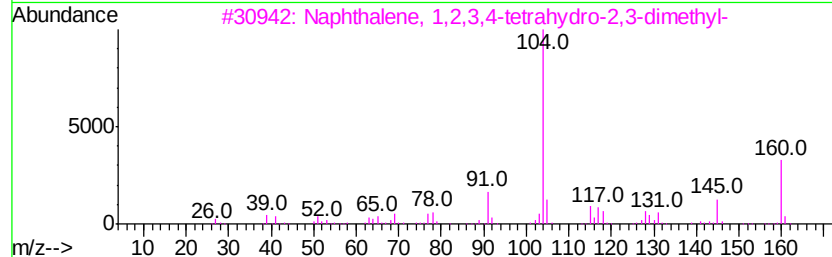
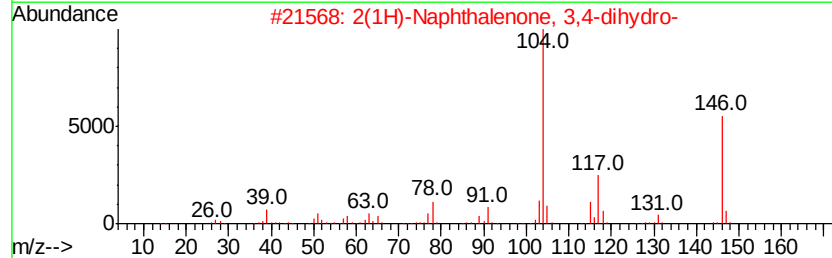
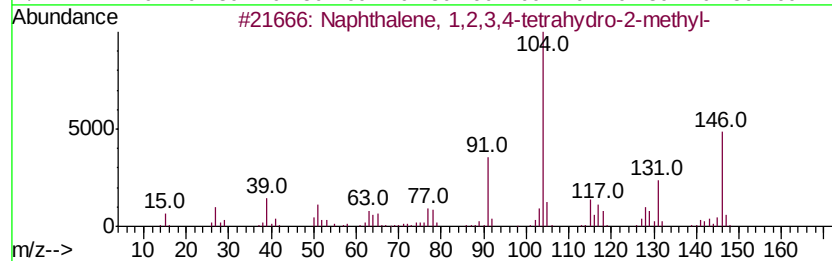
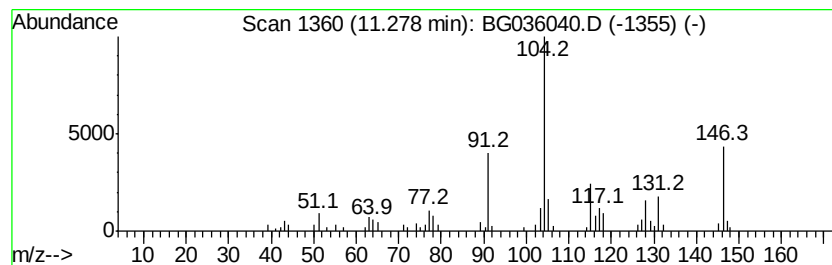
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Naphthalene, 1,2,3,4-tetra... Concentration Rank 15

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.28 | 4.83 ng | 88066 | Naphthalene-d8   | 10.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14    | 003877-19-8  | 87   |
| 2    |    | 2(1H)-Naphthalenone, 3,4-dihydro-   | 146 | C10H10O   | 000530-93-8  | 43   |
| 3    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16    | 021564-92-1  | 42   |
| 4    |    | cis-4-Benzyl-2,6-diphenyltetrahy... | 328 | C24H24O   | 055696-74-7  | 38   |
| 5    |    | 2-Methylbenzyl alcohol, trifluor... | 218 | C10H9F3O2 | 1000364-57-0 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

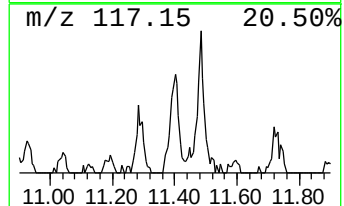
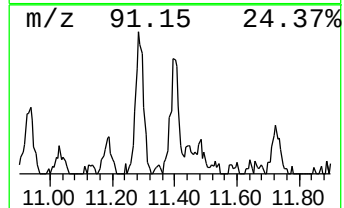
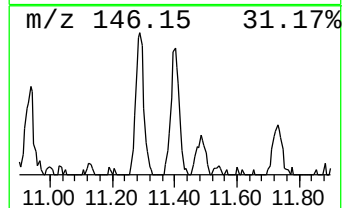
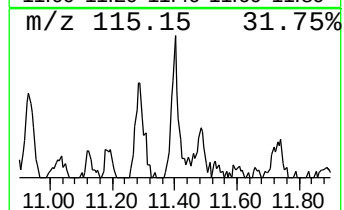
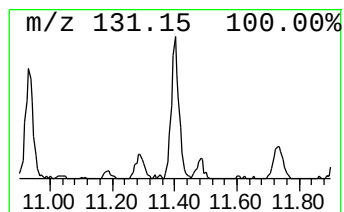
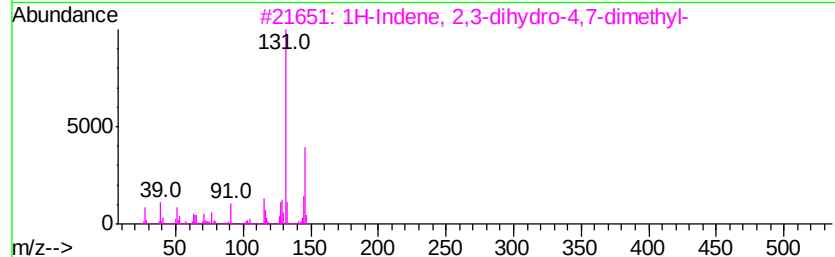
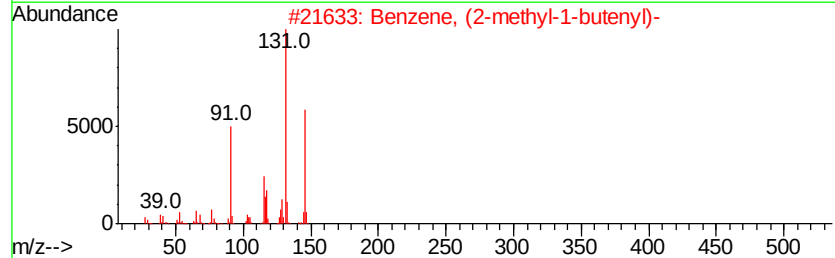
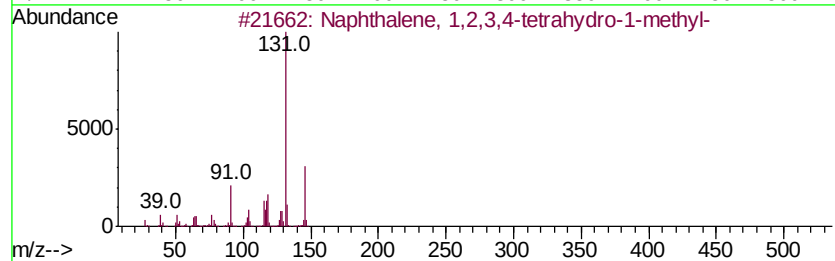
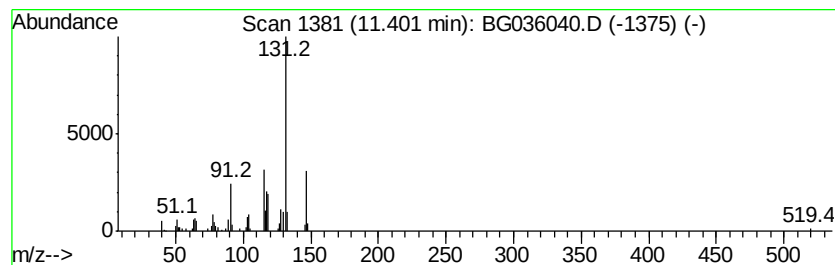
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.40 | 5.10 ng | 93071 | Naphthalene-d8   | 10.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro...  | 146 | C11H14  | 001559-81-5 | 94   |
| 2    |    | Benzene, (2-methyl-1-butenyl)-      | 146 | C11H14  | 056253-64-6 | 90   |
| 3    |    | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 76   |
| 4    |    | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14  | 004175-53-5 | 74   |
| 5    |    | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14  | 017059-48-2 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

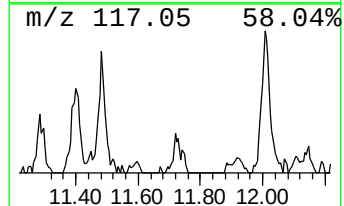
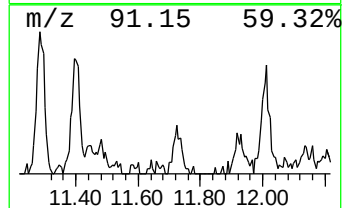
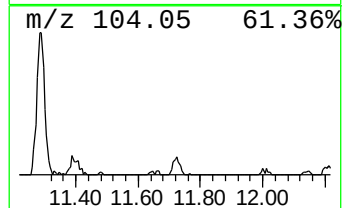
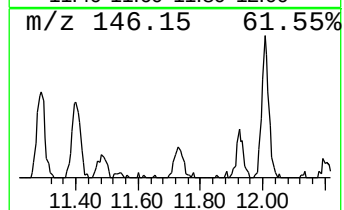
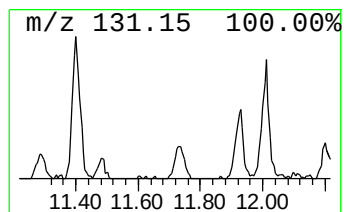
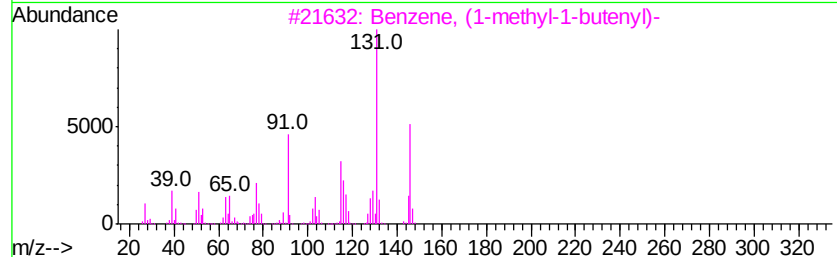
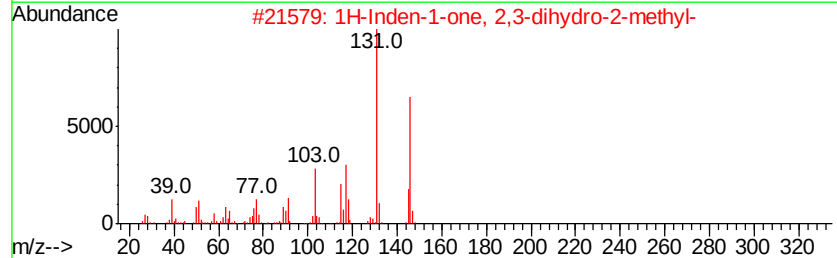
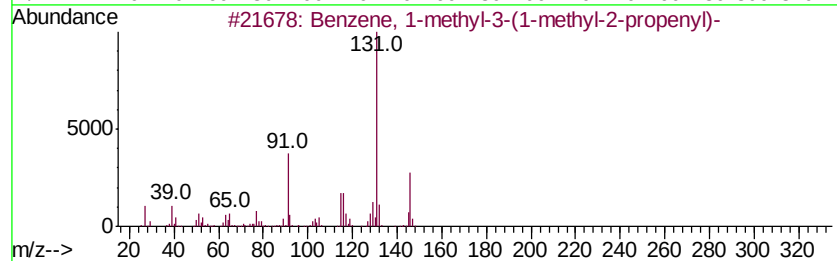
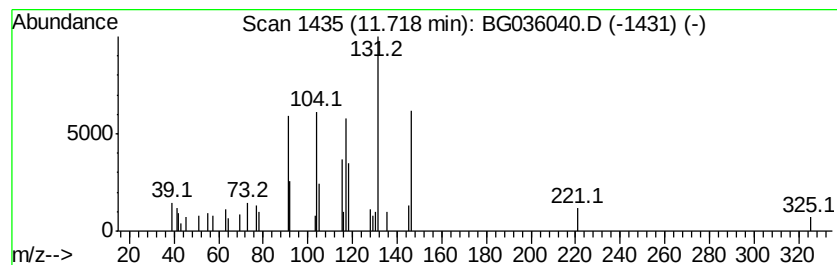
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzene, 1-methyl-3-(1-meth... Concentration Rank 20

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.72 | 2.08 ng | 37886 | Naphthalene-d8   | 10.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-(1-methyl-2-... | 146 | C11H14  | 052161-57-6 | 60   |
| 2    |    | 1H-Inden-1-one, 2,3-dihydro-2-me... | 146 | C10H10O | 017496-14-9 | 58   |
| 3    |    | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 49   |
| 4    |    | Benzene, 1-methyl-2-(1-methyl-2-... | 146 | C11H14  | 097664-19-2 | 49   |
| 5    |    | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 49   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

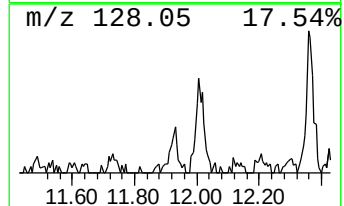
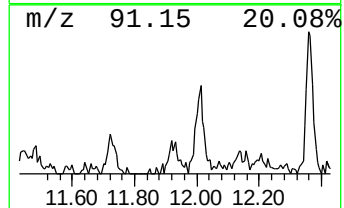
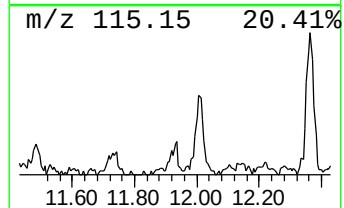
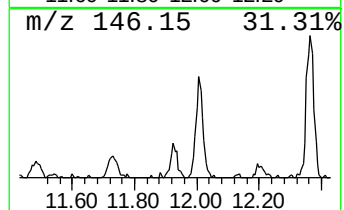
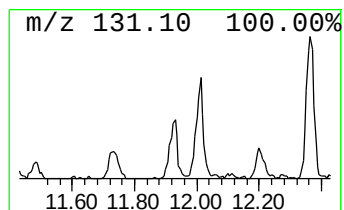
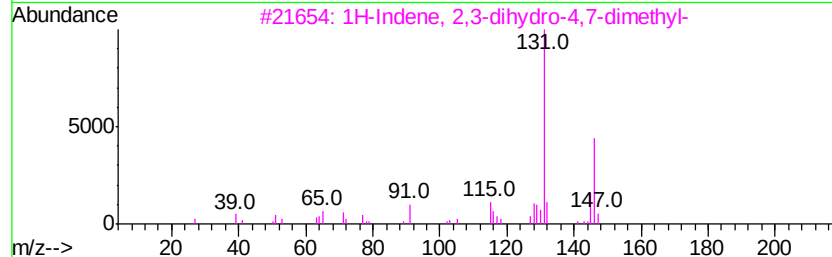
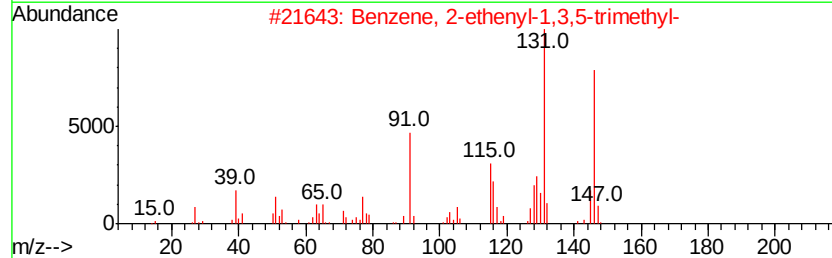
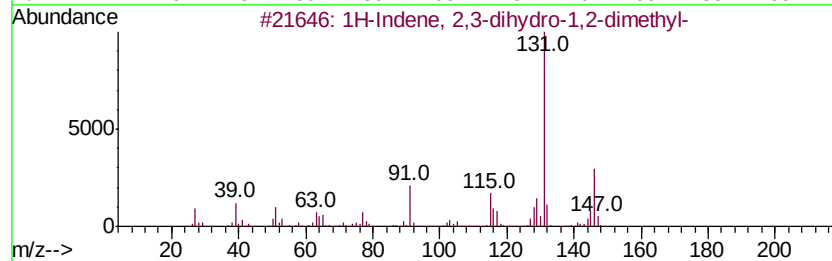
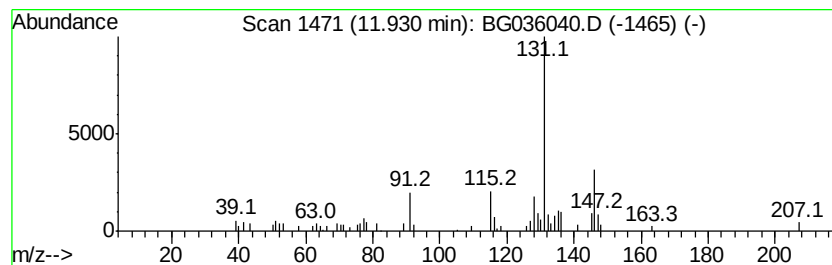
Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 16 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 19

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 11.93   | 2.23 ng                             | 40733        | Napthalene-d8    | 10.83     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 C11H14   | 017057-82-8      | 72        |
| 2       | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 C11H14   | 000769-25-5      | 64        |
| 3       | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 C11H14   | 006682-71-9      | 64        |
| 4       | Benzene, (1,2-dimethyl-1-propenyl)- | 146 C11H14   | 000769-57-3      | 59        |
| 5       | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 C11H14   | 004175-53-5      | 59        |





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

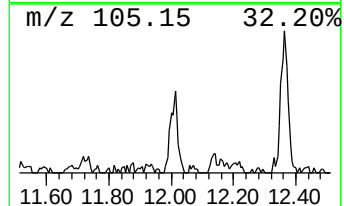
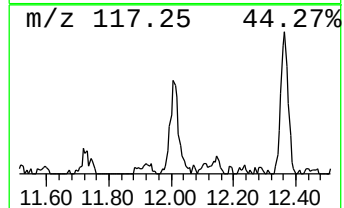
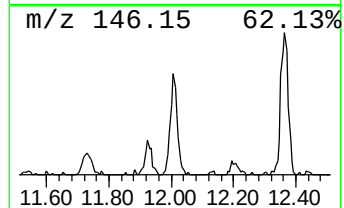
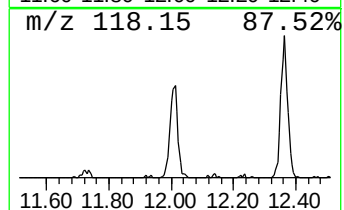
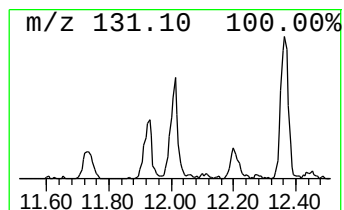
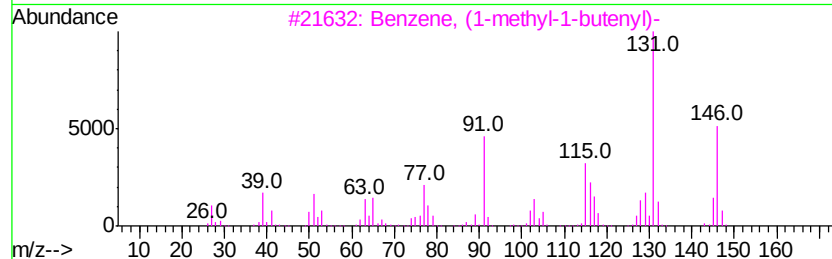
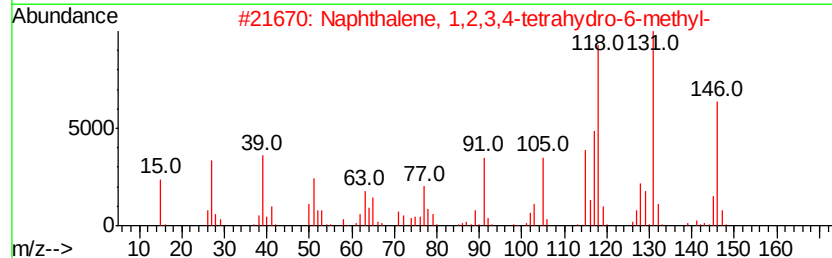
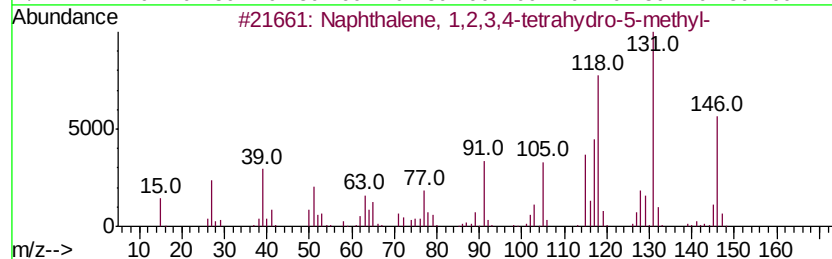
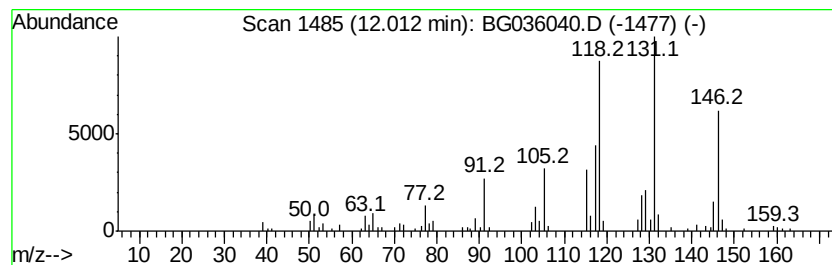
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Naphthalene, 1,2,3,4-tetra... Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.01 | 6.75 ng | 123190 | Naphthalene-d8   | 10.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 002809-64-5 | 93   |
| 2    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001680-51-9 | 90   |
| 3    |    | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 60   |
| 4    |    | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14  | 000769-25-5 | 49   |
| 5    |    | 1H-Inden-1-one, 2,3-dihydro-2-me... | 146 | C10H10O | 017496-14-9 | 45   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

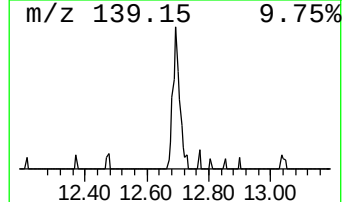
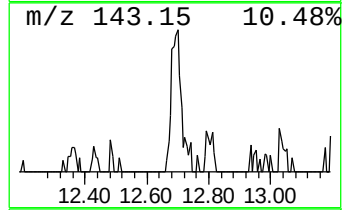
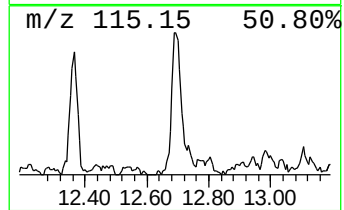
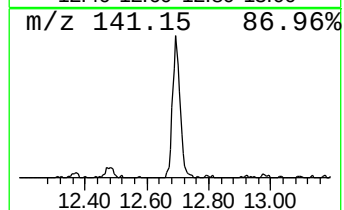
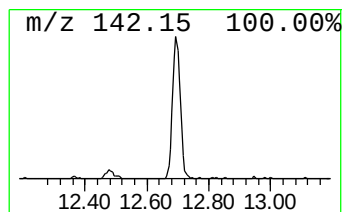
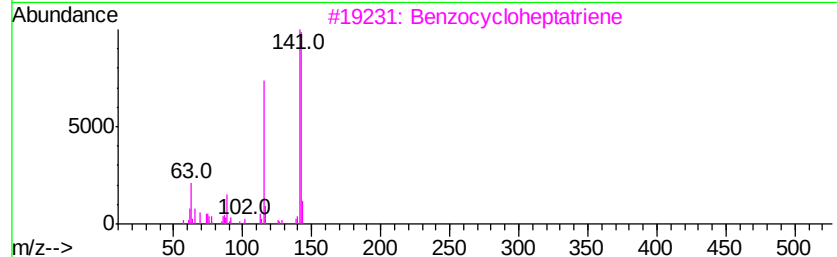
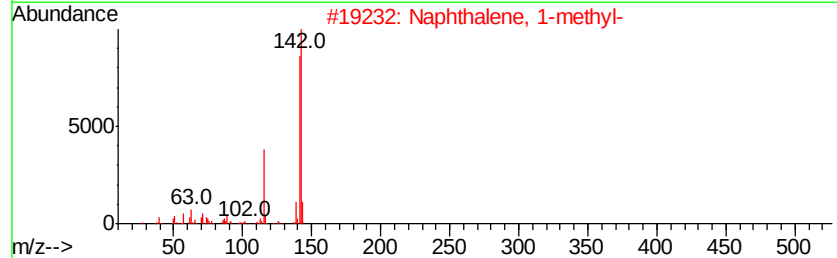
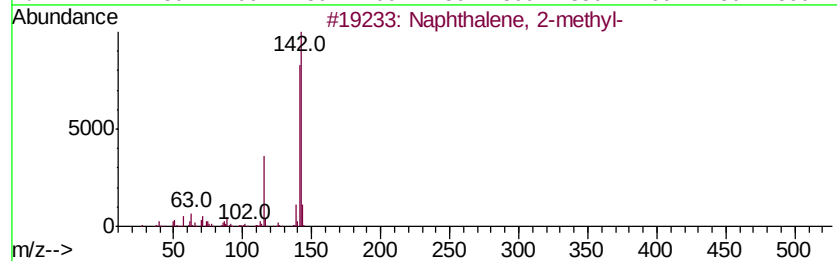
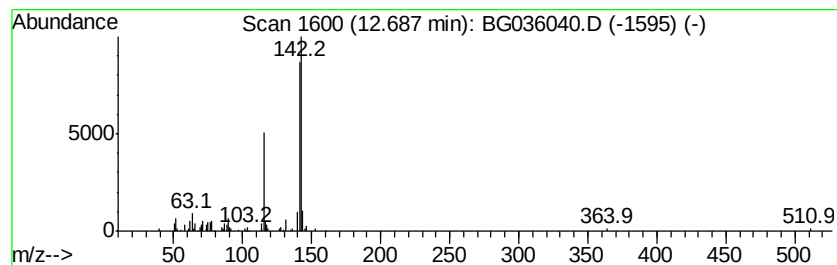
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Naphthalene, 1-methyl- Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.69 | 8.16 ng | 148878 | Naphthalene-d8   | 10.83 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 95   |
| 2    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 95   |
| 3    |    | Benzocycloheptatriene               | 142 | C11H10  | 000264-09-5 | 91   |
| 4    |    | 1H-Indene, 1-ethylidene-            | 142 | C11H10  | 002471-83-2 | 91   |
| 5    |    | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 | C11H10  | 002443-46-1 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080118\  
Data File : BG036040.D  
Acq On : 1 Aug 2018 19:12  
Operator : JU/SJ  
Sample : J4256-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

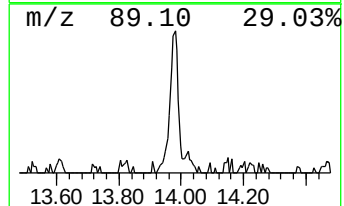
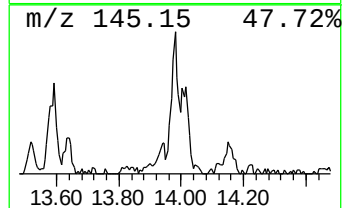
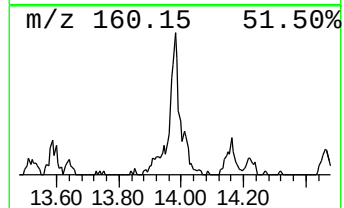
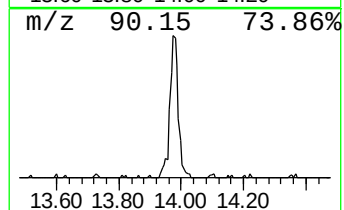
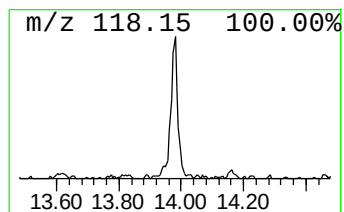
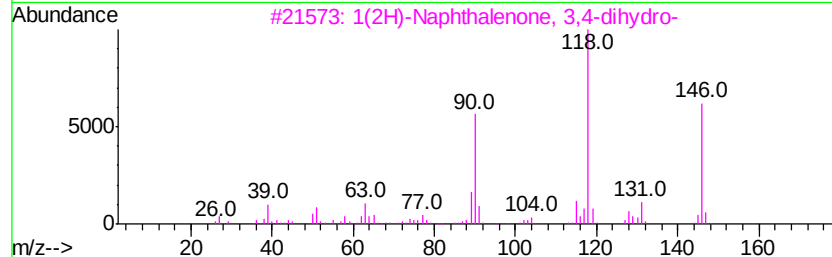
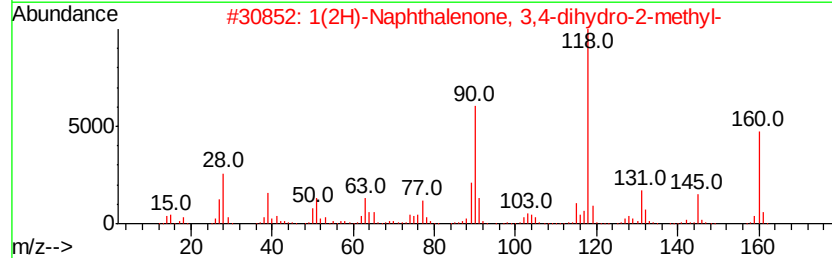
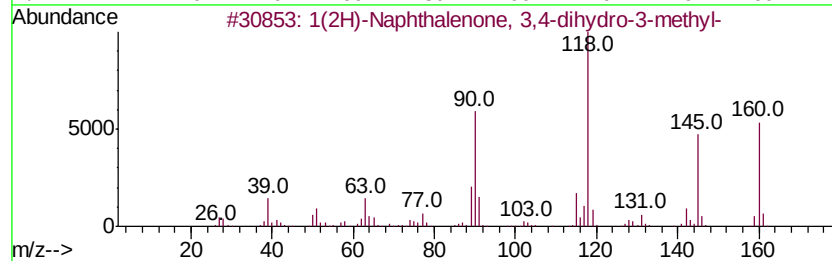
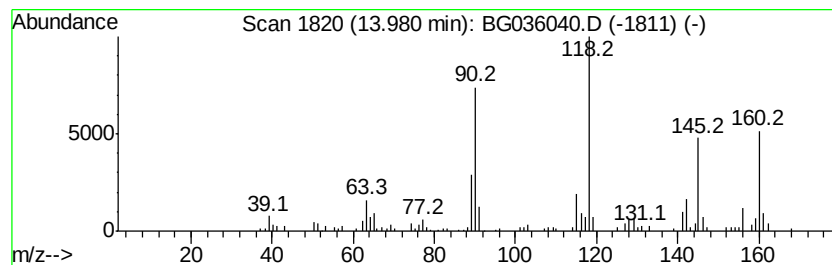
Instrument :  
BNA\_G  
ClientSampleId :  
MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 20 1(2H)-Naphthalenone, 3,4-di... Concentration Rank 12

| R.T.      | EstConc                             | Area   | Relative to ISTD |             | R.T.  |
|-----------|-------------------------------------|--------|------------------|-------------|-------|
| 13.98     | 5.55 ng                             | 156457 | Acenaphthene-d10 |             | 14.64 |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual  |
| 1         | 1(2H)-Naphthalenone, 3,4-dihydro... | 160    | C11H12O          | 014944-23-1 | 94    |
| 2         | 1(2H)-Naphthalenone, 3,4-dihydro... | 160    | C11H12O          | 001590-08-5 | 90    |
| 3         | 1(2H)-Naphthalenone, 3,4-dihydro-   | 146    | C10H10O          | 000529-34-0 | 64    |
| 4         | Homophthalimide                     | 161    | C9H7NO2          | 004456-77-3 | 64    |
| 5         | 1H-2-Benzothiopyran-4(3H)-one       | 164    | C9H8OS           | 004426-76-0 | 59    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG080118\  
 Data File : BG036040.D  
 Acq On : 1 Aug 2018 19:12  
 Operator : JU/SJ  
 Sample : J4256-03  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-A27

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG071218.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 |
| 2-Acetoxyisobutyr... | 5.13  | 5.8     | ng    | 65412    | 1                     | 8.02  | 226370 | 20.0 |
| unknown7.56          | 7.56  | 92.4    | ng    | 1045540  | 1                     | 8.02  | 226370 | 20.0 |
| Indane               | 8.39  | 15.3    | ng    | 173312   | 1                     | 8.02  | 226370 | 20.0 |
| Benzene, 1,4-diet... | 8.51  | 7.6     | ng    | 86002    | 1                     | 8.02  | 226370 | 20.0 |
| Benzene, 1-ethyl-... | 8.66  | 4.0     | ng    | 45287    | 1                     | 8.02  | 226370 | 20.0 |
| Benzene, 2-ethyl-... | 9.13  | 21.9    | ng    | 247326   | 1                     | 8.02  | 226370 | 20.0 |
| o-Cymene             | 9.64  | 6.6     | ng    | 121182   | 2                     | 10.83 | 364756 | 20.0 |
| Benzene, 1,2,3,4-... | 10.22 | 27.1    | ng    | 494679   | 2                     | 10.83 | 364756 | 20.0 |
| 2,2-Dimethylinden... | 10.94 | 3.6     | ng    | 65279    | 2                     | 10.83 | 364756 | 20.0 |
| Naphthalene, 1,2,... | 11.28 | 4.8     | ng    | 88066    | 2                     | 10.83 | 364756 | 20.0 |
| Naphthalene, 1,2,... | 11.40 | 5.1     | ng    | 93071    | 2                     | 10.83 | 364756 | 20.0 |
| Benzene, 1-methyl... | 11.72 | 2.1     | ng    | 37886    | 2                     | 10.83 | 364756 | 20.0 |
| 1H-Indene, 2,3-di... | 11.93 | 2.2     | ng    | 40733    | 2                     | 10.83 | 364756 | 20.0 |
| Naphthalene, 1,2,... | 12.01 | 6.8     | ng    | 123190   | 2                     | 10.83 | 364756 | 20.0 |
| Naphthalene, 1-me... | 12.69 | 8.2     | ng    | 148878   | 2                     | 10.83 | 364756 | 20.0 |
| 1(2H)-Naphthaleno... | 13.98 | 5.5     | ng    | 156457   | 3                     | 14.64 | 563610 | 20.0 |