

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
 Data File : BG053667.D  
 Acq On : 23 May 2022 16:49  
 Operator : CG/JU  
 Sample : N2968-04  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-25

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-BG050522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG053667.D\data.ms

| 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         | 4.222       | 135           | 142         | 158          | rVB      | 2631001        | 4593972       | 57.61%          | 7.263%        |
| 2         | 4.469       | 177           | 184         | 189          | rBV      | 142157         | 219187        | 2.75%           | 0.347%        |
| 3         | 4.533       | 189           | 195         | 205          | rVB      | 140956         | 234037        | 2.94%           | 0.370%        |
| 4         | 5.561       | 361           | 370         | 377          | rBV      | 3168848        | 5481921       | 68.75%          | 8.666%        |
| 5         | 5.632       | 377           | 382         | 386          | rVV      | 249475         | 406943        | 5.10%           | 0.643%        |
| 6         | 5.714       | 386           | 396         | 409          | rVB      | 2859330        | 7973640       | 100.00%         | 12.605%       |
| 7         | 6.067       | 446           | 456         | 466          | rVB      | 2637890        | 4461018       | 55.95%          | 7.052%        |
| 8         | 6.543       | 530           | 537         | 547          | rVB      | 169278         | 283228        | 3.55%           | 0.448%        |
| 9         | 7.030       | 613           | 620         | 628          | rBV      | 407158         | 679112        | 8.52%           | 1.074%        |
| 10        | 7.142       | 632           | 639         | 644          | rVV      | 874619         | 1437869       | 18.03%          | 2.273%        |
| 11        | 7.200       | 644           | 649         | 657          | rVV2     | 628129         | 1255565       | 15.75%          | 1.985%        |
| 12        | 7.277       | 657           | 662         | 674          | rVB      | 561143         | 934724        | 11.72%          | 1.478%        |
| 13        | 7.447       | 683           | 691         | 703          | rBV      | 1025734        | 1705701       | 21.39%          | 2.697%        |
| 14        | 7.712       | 727           | 736         | 744          | rVB      | 3090895        | 5188778       | 65.07%          | 8.203%        |
| 15        | 8.058       | 788           | 795         | 799          | rBV      | 114915         | 205879        | 2.58%           | 0.325%        |
| 16        | 8.176       | 808           | 815         | 823          | rVB      | 1058492        | 1813019       | 22.74%          | 2.866%        |
| 17        | 8.434       | 852           | 859         | 868          | rBV      | 1281425        | 2245627       | 28.16%          | 3.550%        |
| 18        | 8.546       | 868           | 878         | 883          | rVV      | 130760         | 230850        | 2.90%           | 0.365%        |
| 19        | 8.604       | 883           | 888         | 895          | rVB2     | 194106         | 335716        | 4.21%           | 0.531%        |
| 20        | 8.698       | 897           | 904         | 909          | rBV2     | 238866         | 403755        | 5.06%           | 0.638%        |
| 21        | 8.863       | 923           | 932         | 942          | rBV      | 100188         | 187450        | 2.35%           | 0.296%        |
| 22        | 9.016       | 951           | 958         | 962          | rBV      | 215952         | 368399        | 4.62%           | 0.582%        |
| 23        | 9.063       | 962           | 966         | 973          | rVV      | 161850         | 273032        | 3.42%           | 0.432%        |
| 24        | 9.168       | 975           | 984         | 989          | rVV      | 507995         | 888285        | 11.14%          | 1.404%        |
| 25        | 9.227       | 989           | 994         | 1009         | rVB2     | 485564         | 1427358       | 17.90%          | 2.256%        |
| 26        | 9.503       | 1033          | 1041        | 1047         | rBV2     | 106692         | 195471        | 2.45%           | 0.309%        |
| 27        | 9.691       | 1066          | 1073        | 1078         | rBV      | 280240         | 464641        | 5.83%           | 0.735%        |
| 28        | 9.750       | 1078          | 1083        | 1092         | rVB      | 347196         | 581730        | 7.30%           | 0.920%        |
| 29        | 10.114      | 1139          | 1145        | 1155         | rVB      | 388526         | 701836        | 8.80%           | 1.110%        |
| 30        | 10.267      | 1158          | 1171        | 1179         | rBV2     | 858789         | 1699369       | 21.31%          | 2.686%        |
| 31        | 10.502      | 1206          | 1211        | 1217         | rVB      | 111254         | 190156        | 2.38%           | 0.301%        |
| 32        | 10.866      | 1262          | 1273        | 1277         | rVV2     | 174566         | 521644        | 6.54%           | 0.825%        |
| 33        | 10.925      | 1277          | 1283        | 1289         | rVV      | 1771039        | 3288320       | 41.24%          | 5.198%        |
| 34        | 10.984      | 1289          | 1293        | 1297         | rVB      | 79204          | 121969        | 1.53%           | 0.193%        |
| 35        | 11.042      | 1297          | 1303        | 1308         | rBV      | 165331         | 293456        | 3.68%           | 0.464%        |
| 36        | 11.530      | 1381          | 1386        | 1391         | rVB      | 66683          | 121905        | 1.53%           | 0.193%        |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
 Data File : BG053667.D  
 Acq On : 23 May 2022 16:49  
 Operator : CG/JU  
 Sample : N2968-04  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-25

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-BG050522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 11.782 | 1423 | 1429 | 1436 | rVB  | 91784   | 162091  | 2.03%  | 0.256% |
| 38 | 11.970 | 1454 | 1461 | 1471 | rVV2 | 106133  | 233297  | 2.93%  | 0.369% |
| 39 | 12.053 | 1471 | 1475 | 1490 | rVV4 | 53482   | 130267  | 1.63%  | 0.206% |
| 40 | 12.247 | 1502 | 1508 | 1517 | rVB3 | 90600   | 167797  | 2.10%  | 0.265% |
| 41 | 12.417 | 1529 | 1537 | 1543 | rVB3 | 76406   | 138684  | 1.74%  | 0.219% |
| 42 | 12.523 | 1543 | 1555 | 1561 | rBV  | 324583  | 582723  | 7.31%  | 0.921% |
| 43 | 12.740 | 1581 | 1592 | 1603 | rVV  | 586276  | 1080464 | 13.55% | 1.708% |
| 44 | 13.151 | 1656 | 1662 | 1671 | rVB4 | 50519   | 93630   | 1.17%  | 0.148% |
| 45 | 13.310 | 1680 | 1689 | 1696 | rBV  | 1231858 | 1922834 | 24.11% | 3.040% |
| 46 | 13.856 | 1778 | 1782 | 1791 | rVB2 | 128626  | 255774  | 3.21%  | 0.404% |
| 47 | 14.009 | 1801 | 1808 | 1813 | rBV  | 110710  | 211904  | 2.66%  | 0.335% |
| 48 | 14.062 | 1813 | 1817 | 1825 | rVB4 | 84359   | 148392  | 1.86%  | 0.235% |
| 49 | 14.690 | 1913 | 1924 | 1932 | rBV  | 301010  | 552934  | 6.93%  | 0.874% |
| 50 | 16.177 | 2169 | 2177 | 2187 | rBV  | 1009490 | 1561070 | 19.58% | 2.468% |
| 51 | 17.434 | 2385 | 2391 | 2396 | rBV  | 370200  | 558925  | 7.01%  | 0.884% |
| 52 | 18.321 | 2537 | 2542 | 2546 | rBV2 | 51114   | 88977   | 1.12%  | 0.141% |
| 53 | 20.042 | 2828 | 2835 | 2841 | rBV  | 2075034 | 2781677 | 34.89% | 4.397% |
| 54 | 21.710 | 3113 | 3119 | 3127 | rVB  | 360011  | 590736  | 7.41%  | 0.934% |
| 55 | 24.970 | 3665 | 3674 | 3685 | rVB2 | 202416  | 578214  | 7.25%  | 0.914% |

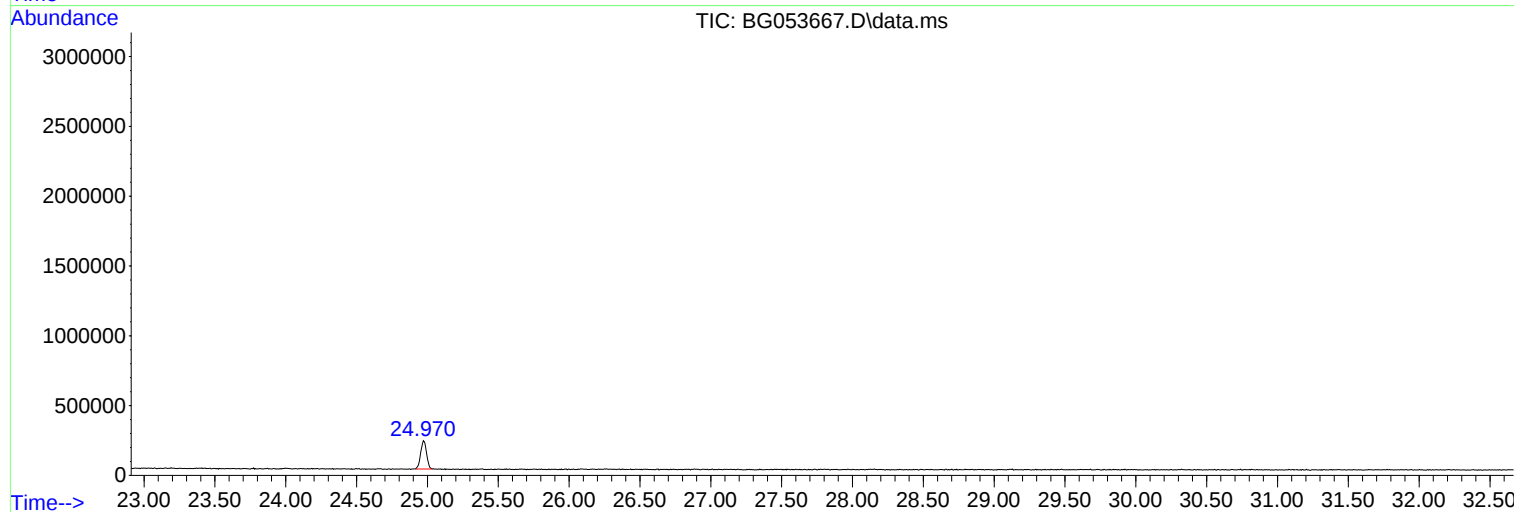
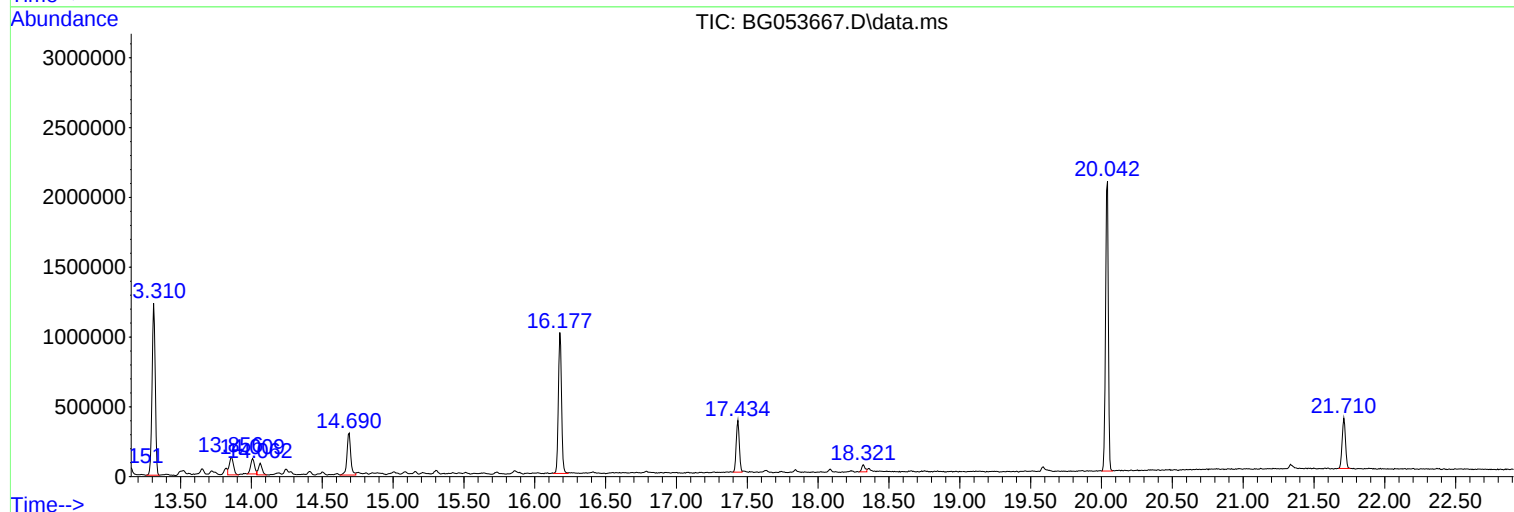
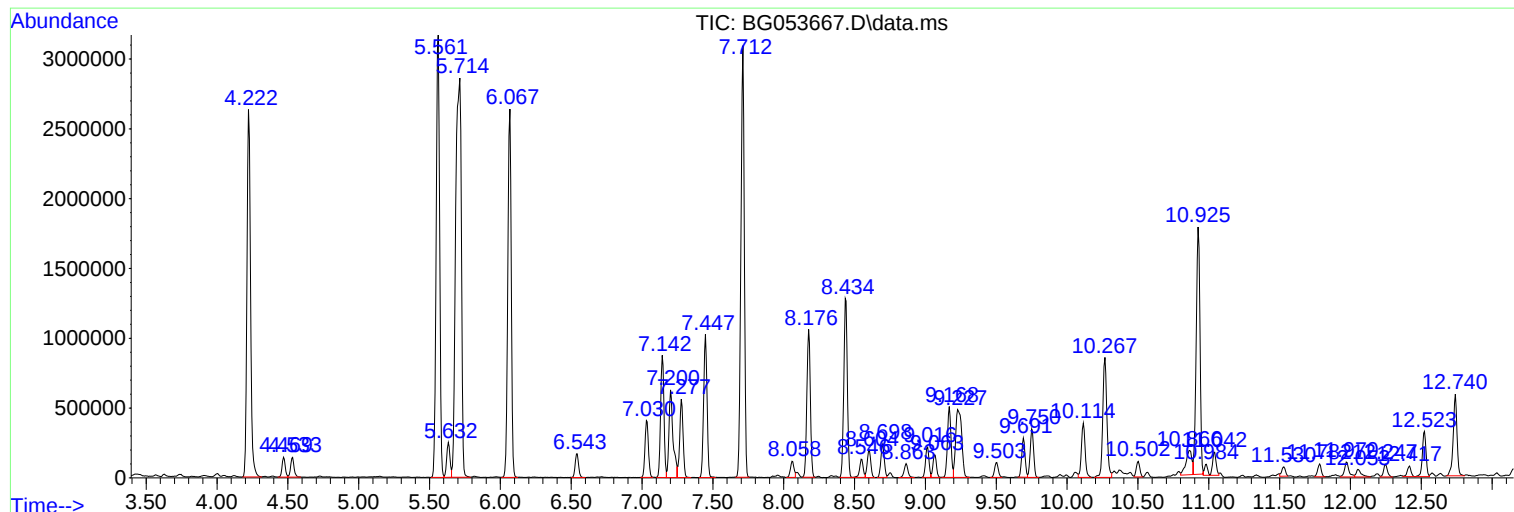
Sum of corrected areas: 63255952

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

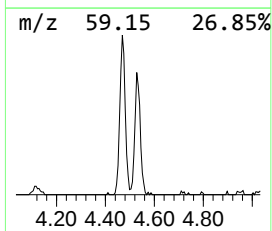
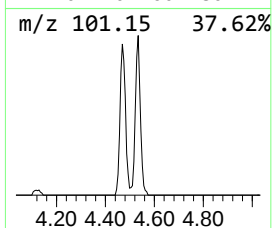
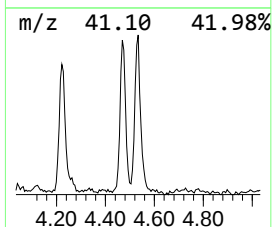
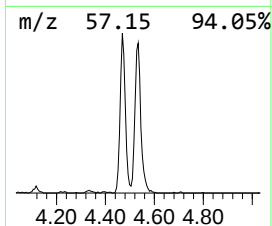
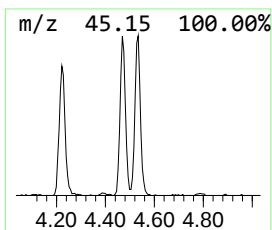
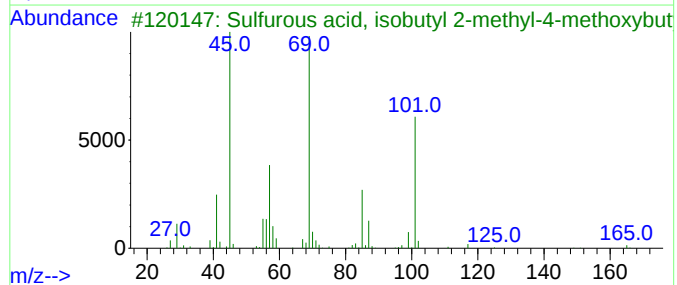
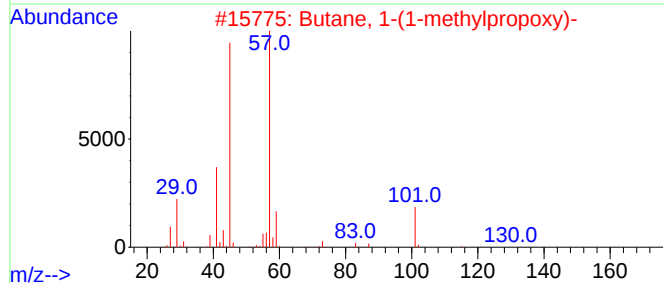
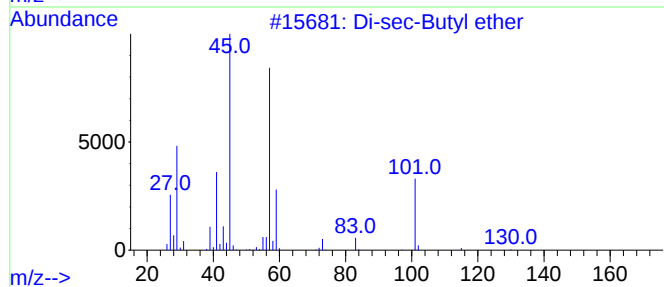
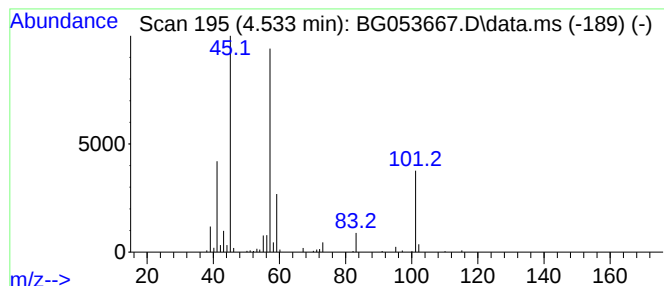
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Di-sec-Butyl ether Concentration Rank 19

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 4.533 | 22.74 ng | 234037 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Di-sec-Butyl ether                  | 130 | C8H18O    | 006863-58-7  | 90   |
| 2    |    |   | Butane, 1-(1-methylpropoxy)-        | 130 | C8H18O    | 000999-65-5  | 78   |
| 3    |    |   | Sulfurous acid, isobutyl 2-methy... | 238 | C10H22O4S | 1000309-20-3 | 42   |
| 4    |    |   | 3-Pentanol, 3-ethyl-2-methyl-       | 130 | C8H18O    | 000597-05-7  | 39   |
| 5    |    |   | 2-Hydroxy-3-pentanone               | 102 | C5H10O2   | 005704-20-1  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

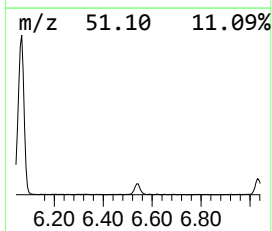
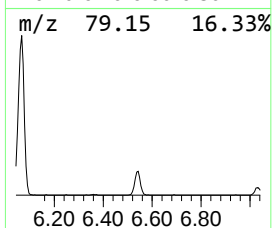
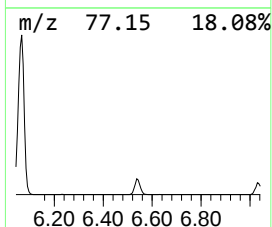
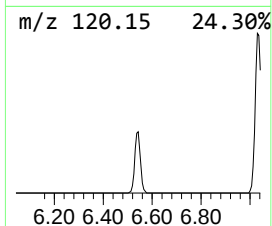
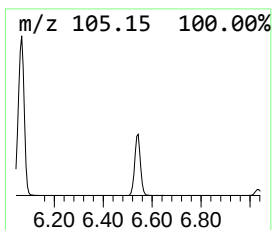
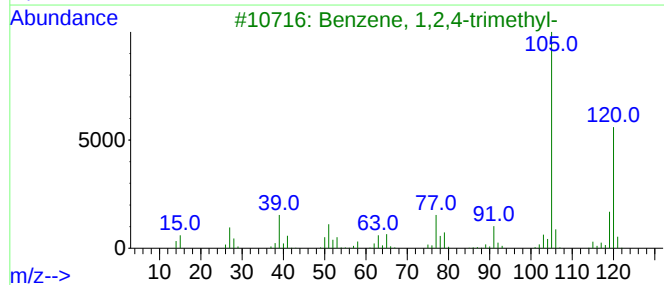
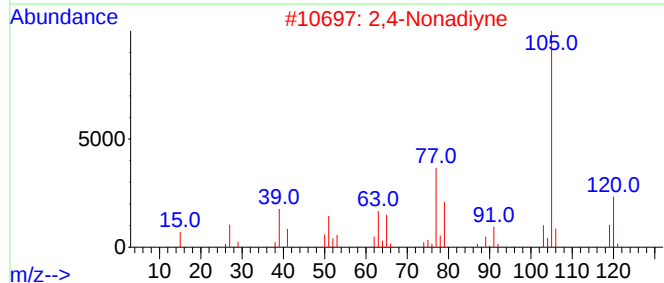
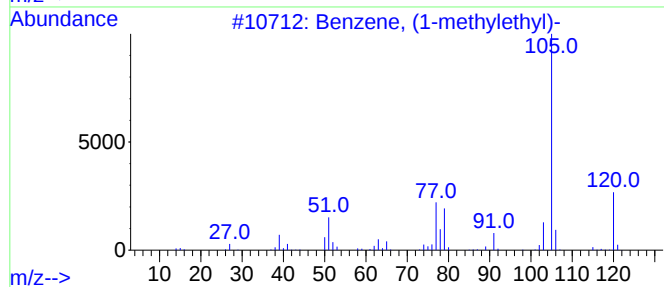
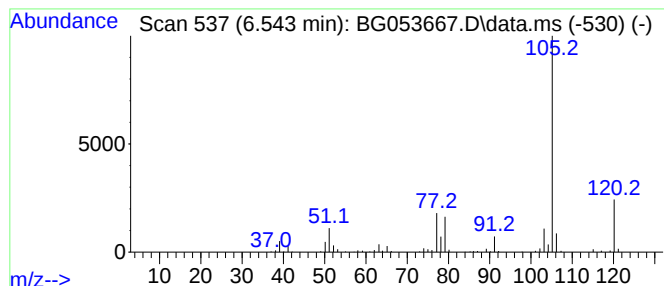
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Benzene, (1-methylethyl)- Concentration Rank 16

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 6.543 | 27.51 ng | 283228 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methylethyl)-  | 120 | C9H12   | 000098-82-8 | 93   |
| 2    |    |   | 2,4-Nonadiyne              | 120 | C9H12   | 063621-15-8 | 80   |
| 3    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 78   |
| 4    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 78   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

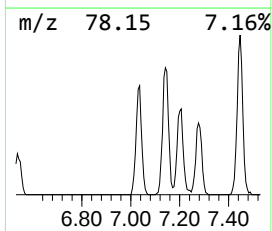
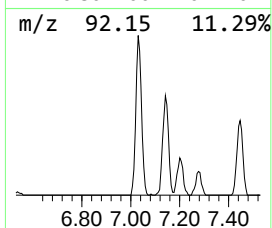
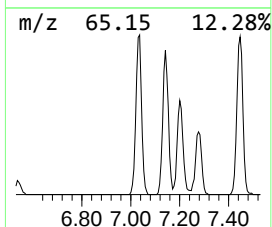
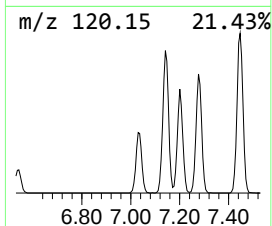
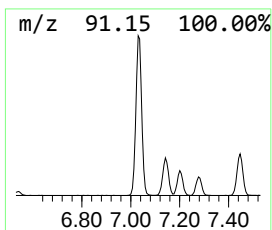
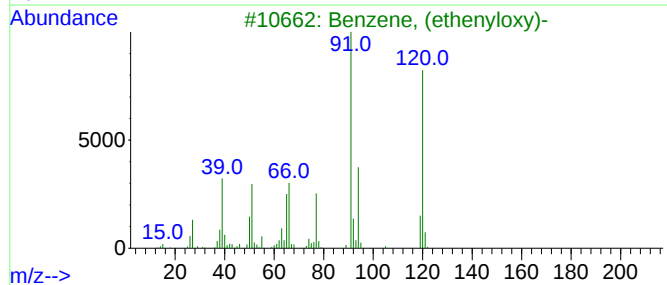
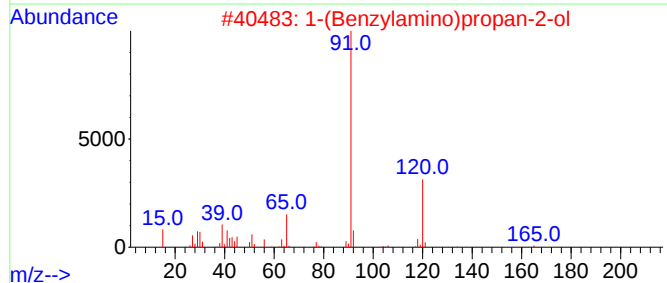
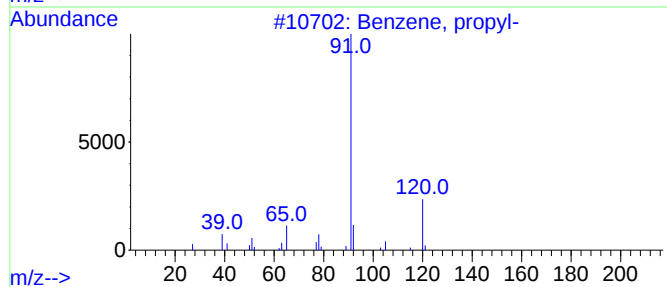
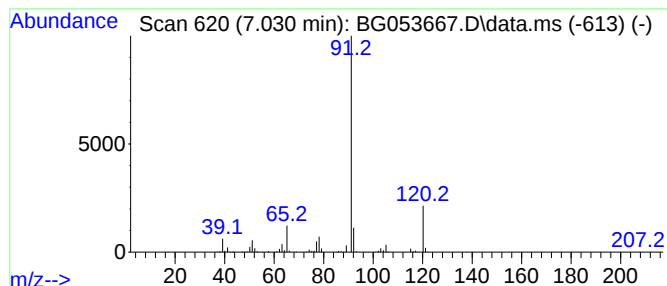
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Benzene, propyl- Concentration Rank 11

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 7.030 | 65.97 ng | 679112 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Benzene, propyl-                    | 120 | C9H12     | 000103-65-1  | 94   |
| 2    |    |   | 1-(Benzylamino)propan-2-ol          | 165 | C10H15NO  | 1000486-84-7 | 72   |
| 3    |    |   | Benzene, (ethenyloxy)-              | 120 | C8H8O     | 000766-94-9  | 64   |
| 4    |    |   | Hydrazinecarboxylic acid, phenyl... | 166 | C8H10N2O2 | 005331-43-1  | 50   |
| 5    |    |   | Benzeneacetaldehyde                 | 120 | C8H8O     | 000122-78-1  | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

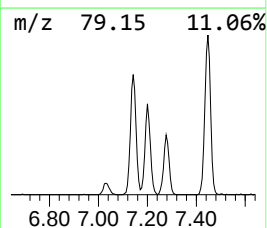
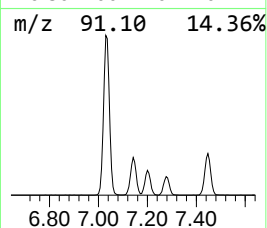
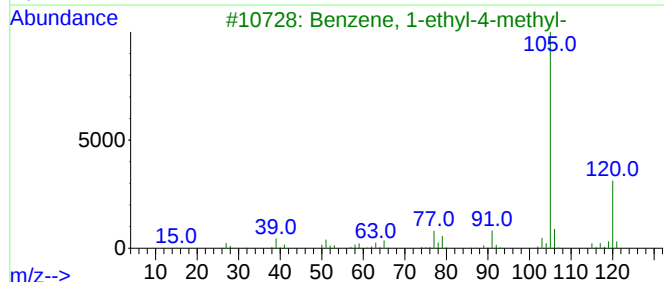
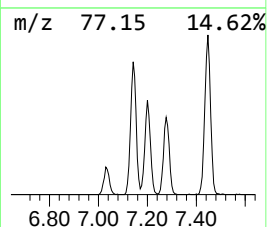
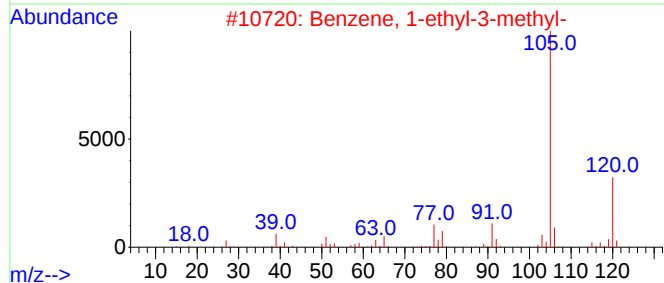
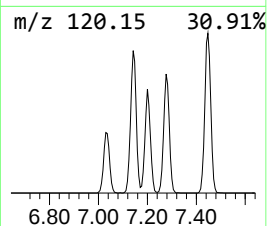
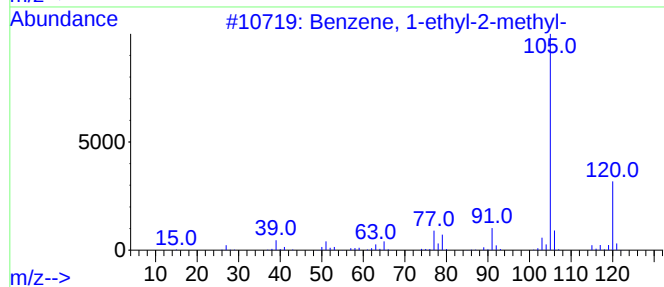
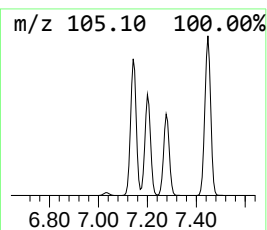
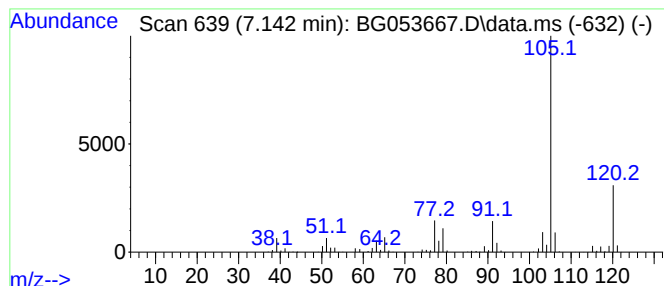
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 7.142 | 139.68 ng | 1437870 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 3    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 4    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

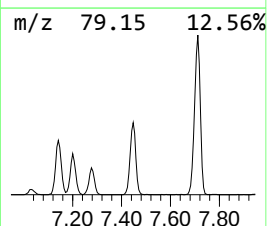
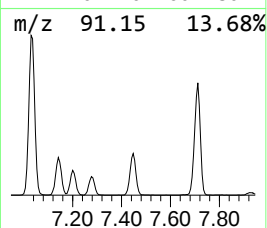
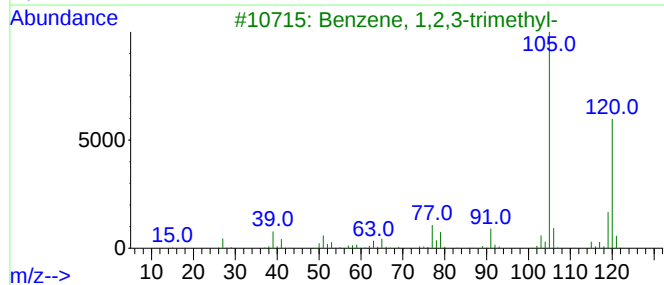
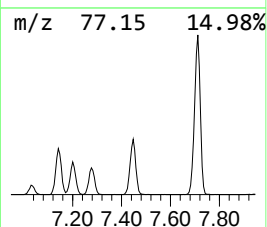
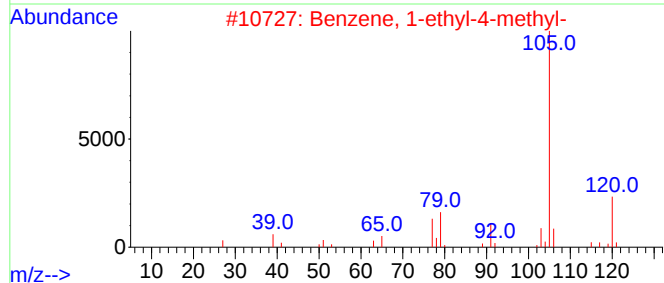
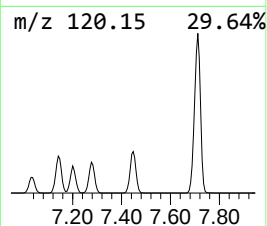
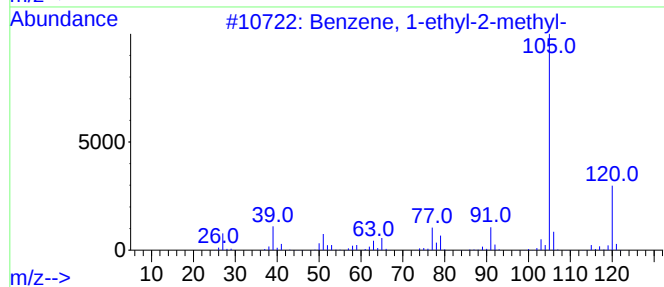
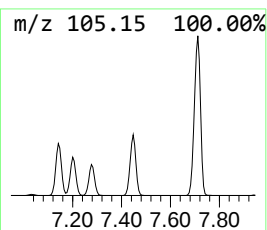
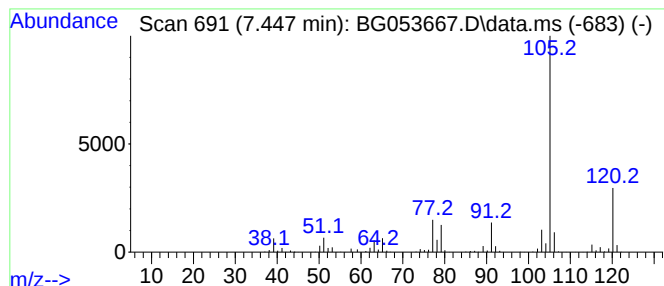
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Benzene, 1-ethyl-4-methyl- Concentration Rank 8

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 7.447 | 165.70 ng | 1705700 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 2    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 93   |
| 3    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 4    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

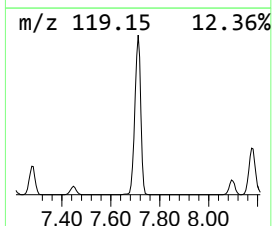
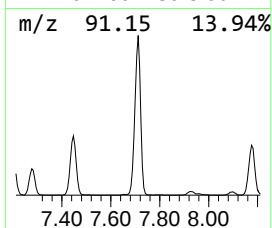
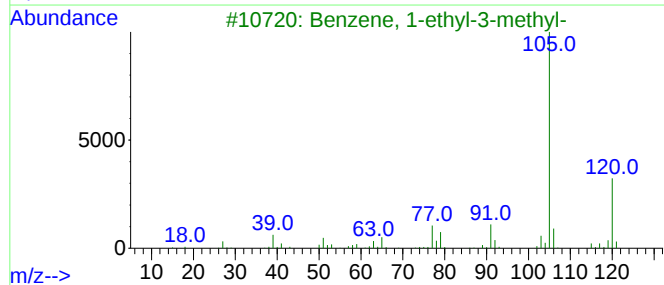
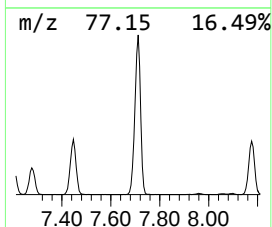
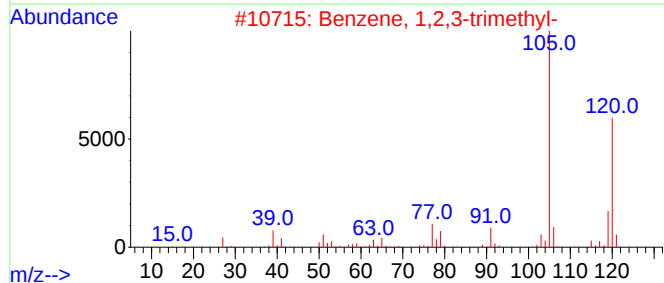
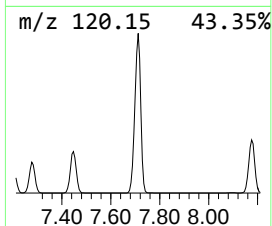
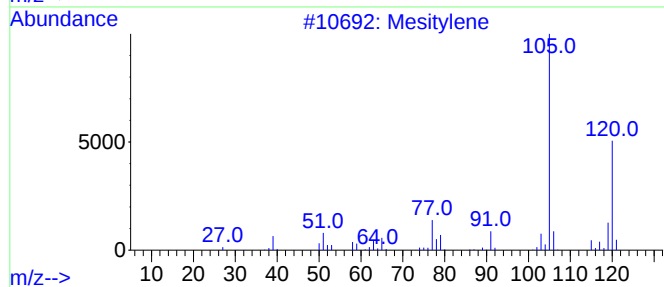
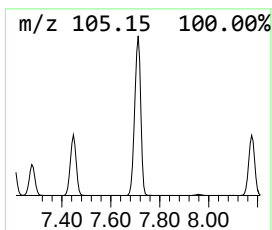
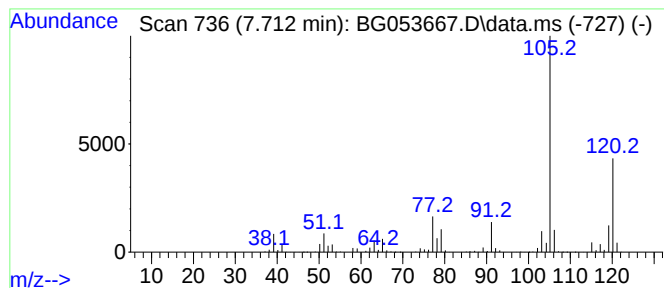
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Mesitylene Concentration Rank 3

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 7.712 | 504.06 ng | 5188780 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 97   |
| 2    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 96   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

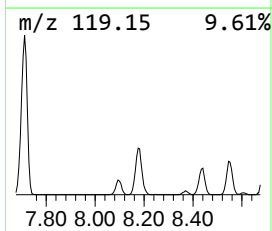
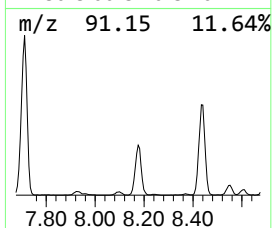
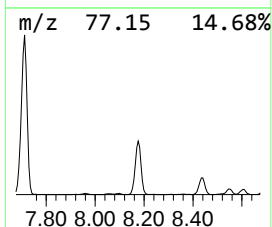
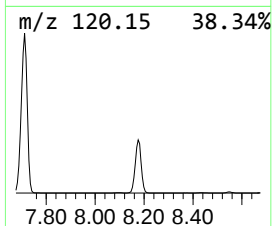
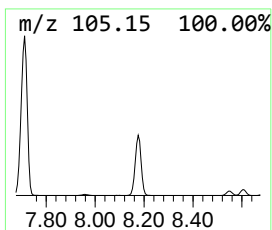
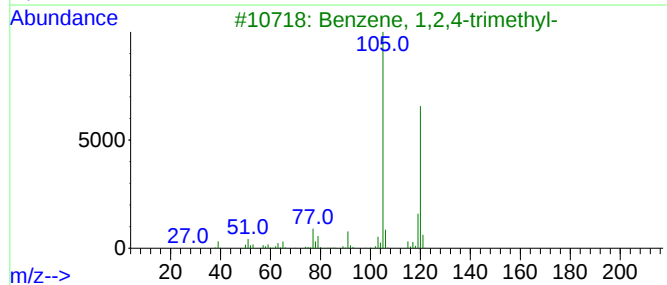
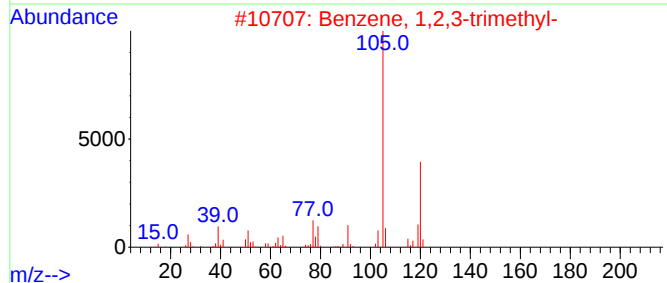
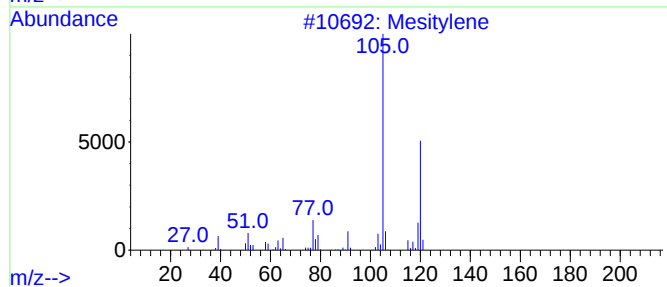
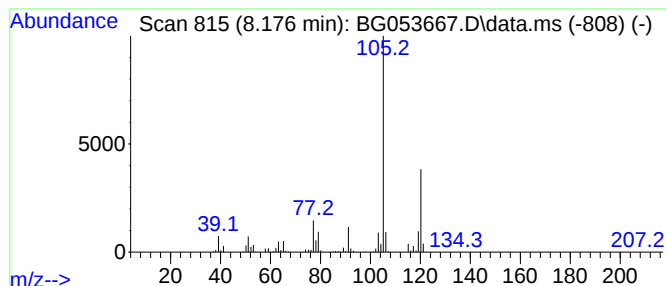
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 7

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 8.176 | 176.13 ng | 1813020 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 95   |
| 2    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 95   |
| 3    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |
| 5    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

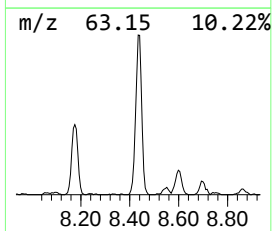
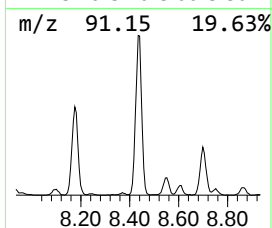
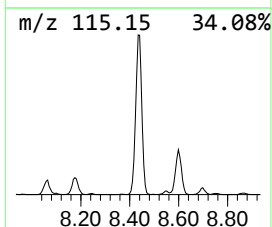
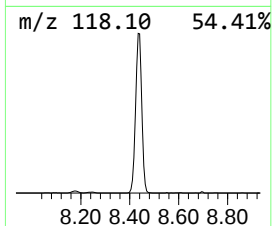
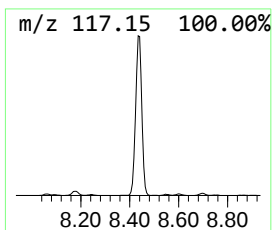
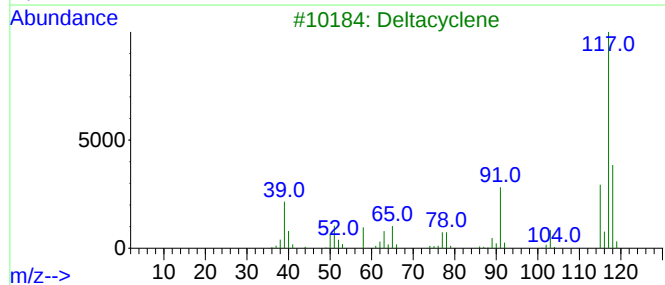
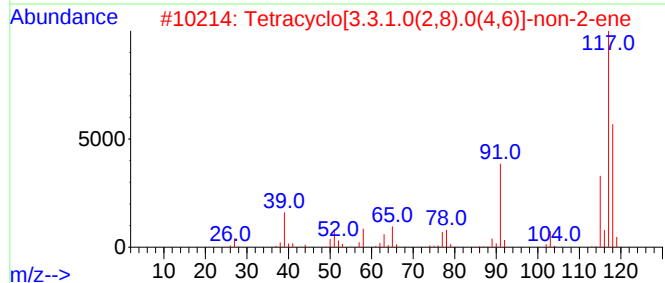
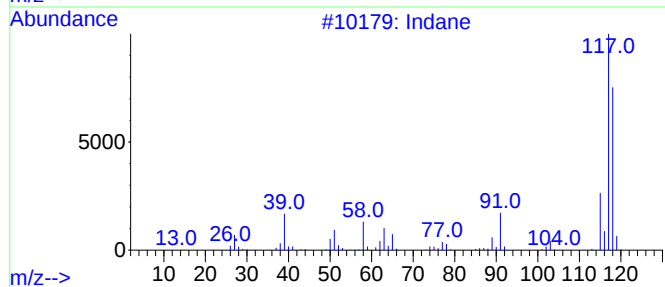
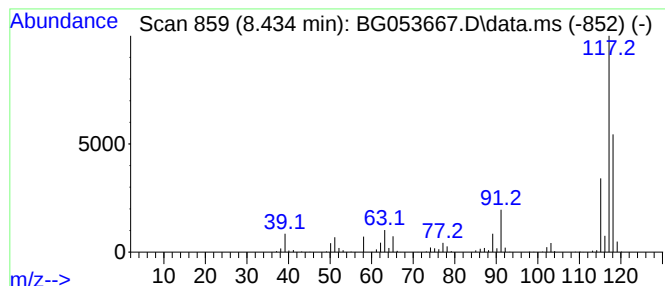
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Indane Concentration Rank 6

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 8.434 | 218.15 ng | 2245630 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|------|-------------------------------------|-----|---------|--------------|------|
| 1    |      | Indane                              | 118 | C9H10   | 000496-11-7  | 91   |
| 2    |      | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10   | 1000191-13-7 | 72   |
| 3    |      | Deltacyclene                        | 118 | C9H10   | 007785-10-6  | 68   |
| 4    |      | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4  | 68   |
| 5    |      | Benzene, 1-ethenyl-2-methyl-        | 118 | C9H10   | 000611-15-4  | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

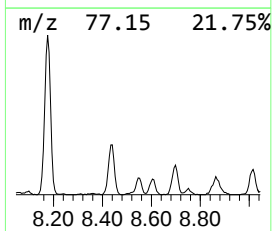
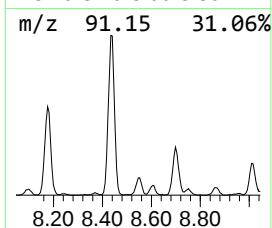
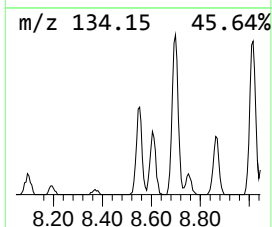
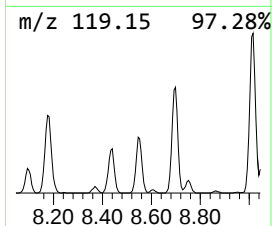
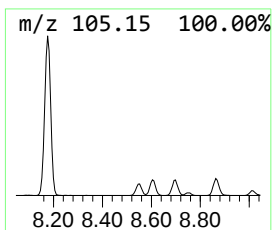
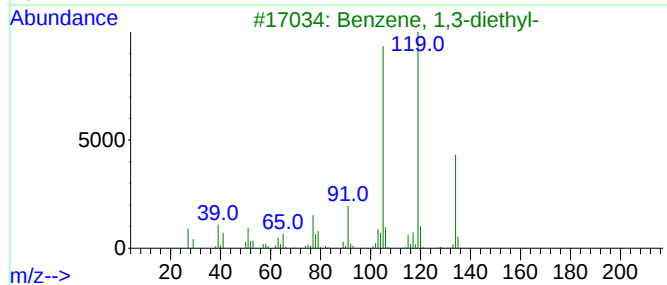
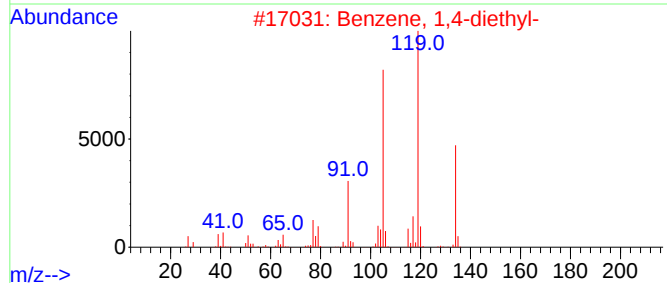
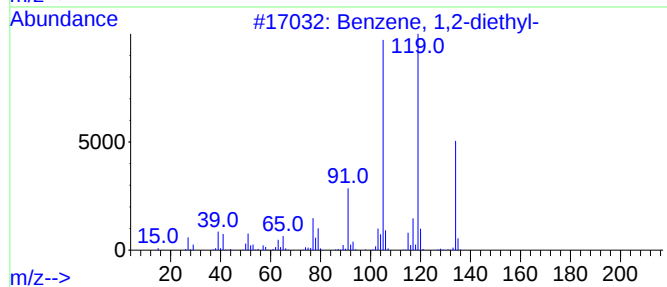
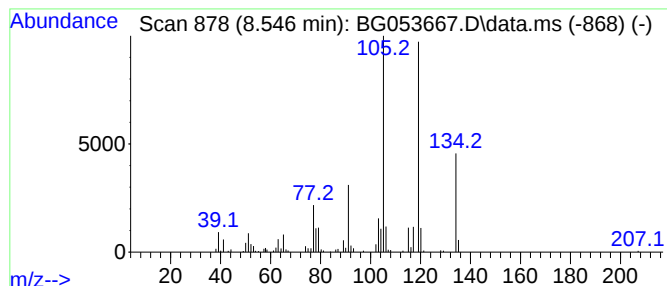
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Benzene, 1,2-diethyl- Concentration Rank 20

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 8.546 | 22.43 ng | 230850 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 95   |
| 2    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 94   |
| 3    |    |   | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 94   |
| 4    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 87   |
| 5    |    |   | 1,3,8-p-Menthatriene           | 134 | C10H14  | 018368-95-1 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

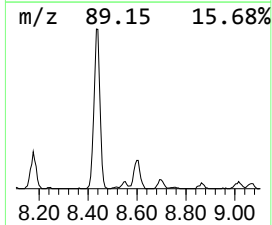
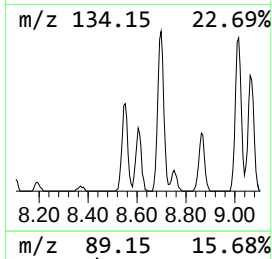
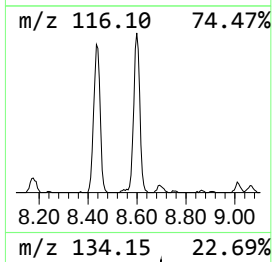
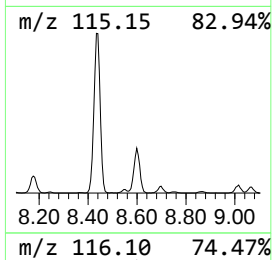
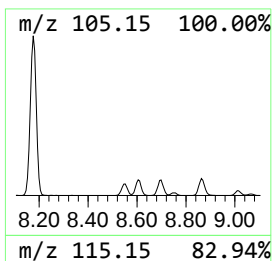
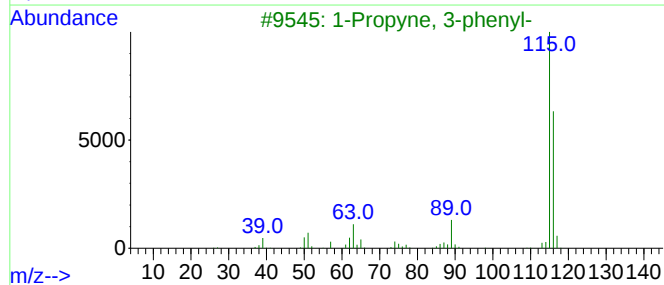
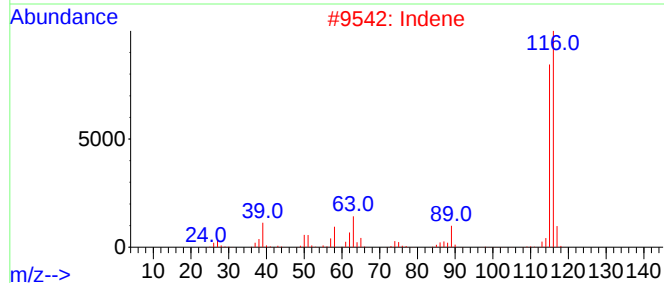
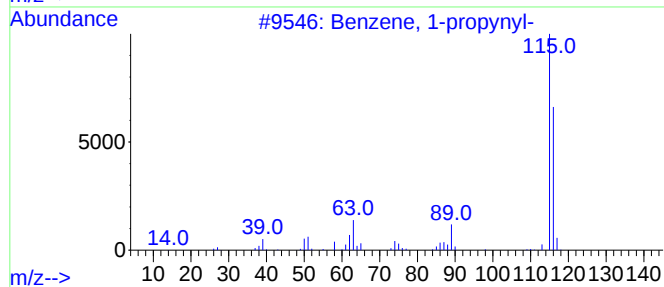
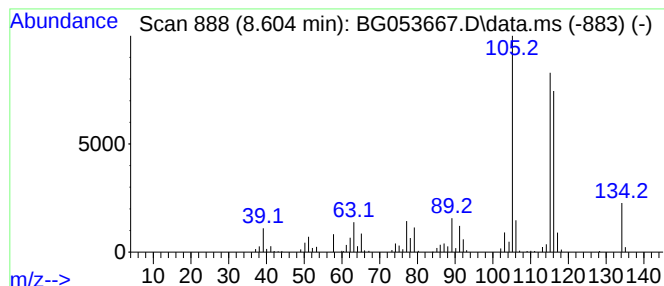
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Benzene, 1-propynyl- Concentration Rank 15

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 8.604 | 32.61 ng | 335716 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-propynyl-         | 116 | C9H8    | 000673-32-5 | 93   |
| 2    |    |   | Indene                       | 116 | C9H8    | 000095-13-6 | 93   |
| 3    |    |   | 1-Propyne, 3-phenyl-         | 116 | C9H8    | 010147-11-2 | 92   |
| 4    |    |   | 3-Methylphenylacetylene      | 116 | C9H8    | 000766-82-5 | 92   |
| 5    |    |   | Benzene, 1-ethynyl-4-methyl- | 116 | C9H8    | 000766-97-2 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

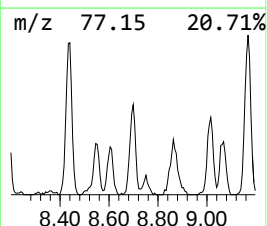
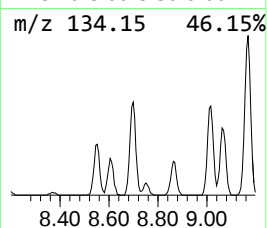
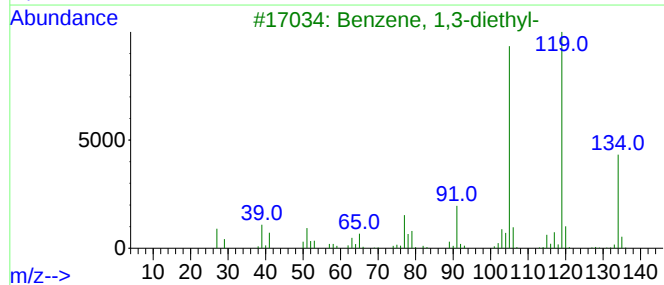
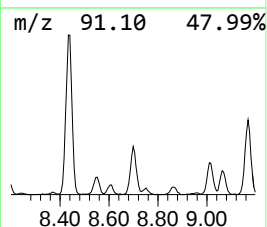
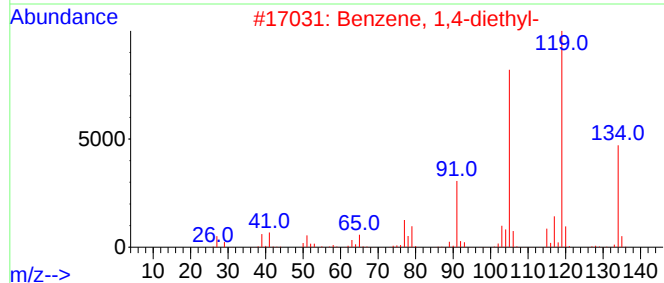
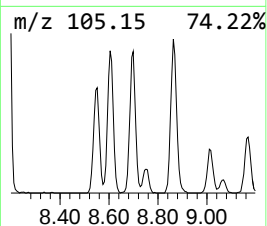
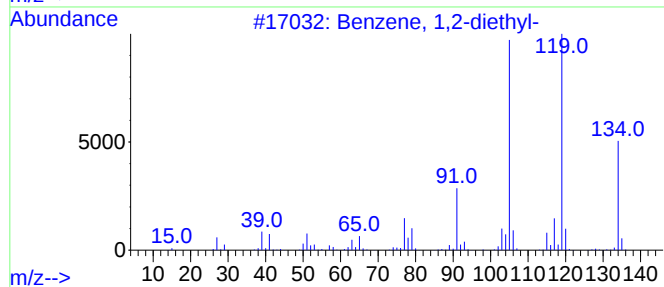
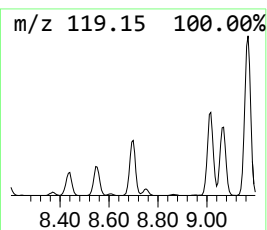
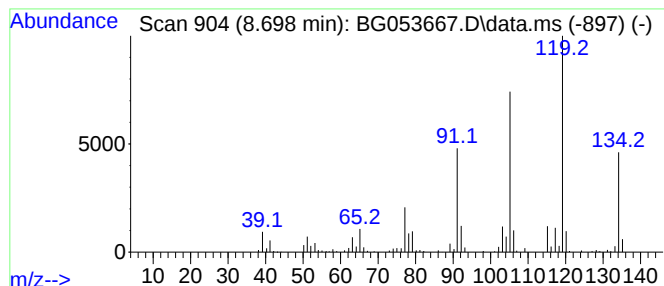
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzene, 1,4-diethyl- Concentration Rank 13

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 8.698 | 39.22 ng | 403755 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 95   |
| 2    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 94   |
| 3    |    |   | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 94   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 90   |
| 5    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

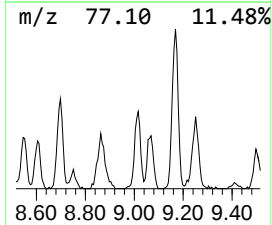
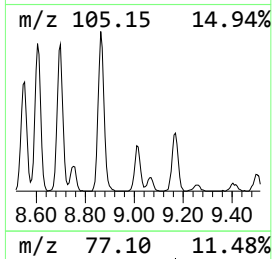
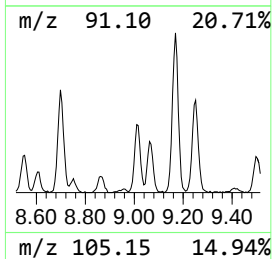
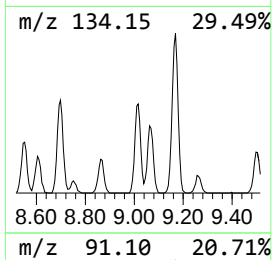
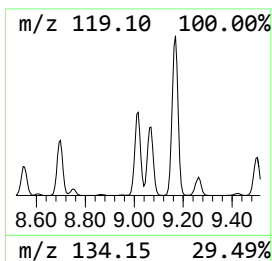
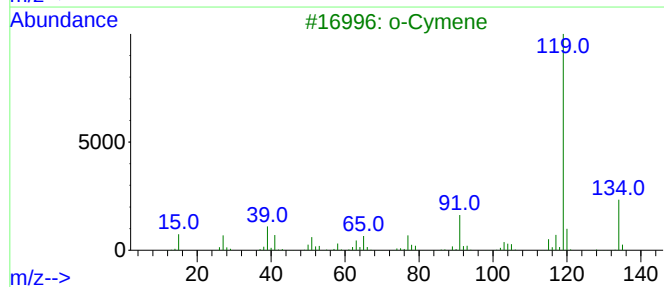
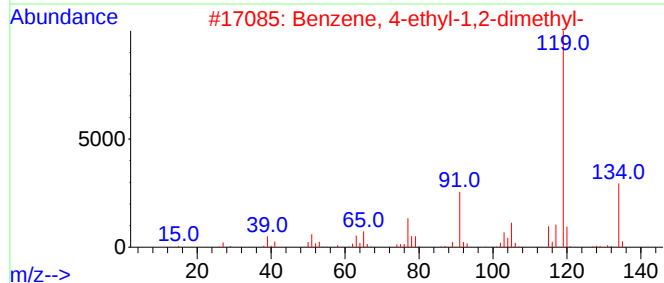
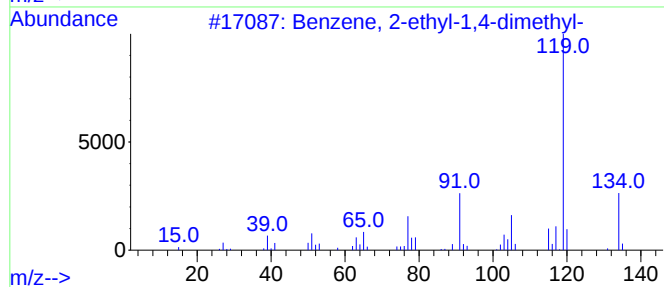
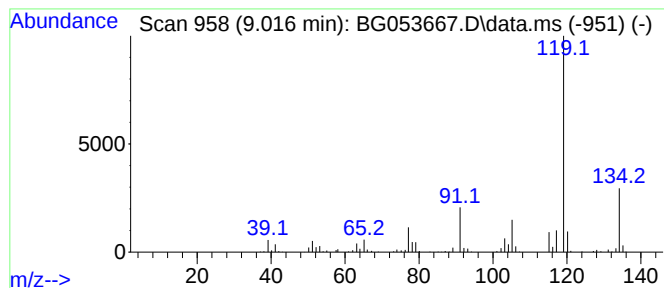
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 14

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 9.016 | 35.79 ng | 368399 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 96   |
| 2    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 96   |
| 3    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 94   |
| 4    |    |   | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 94   |
| 5    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

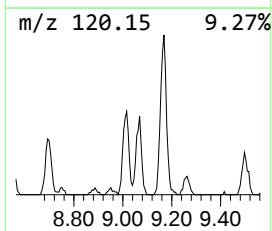
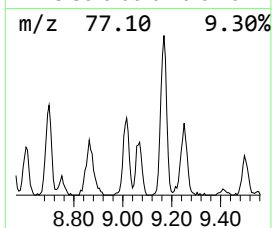
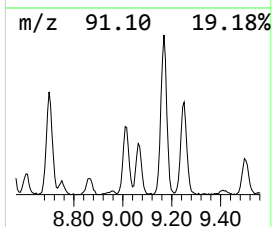
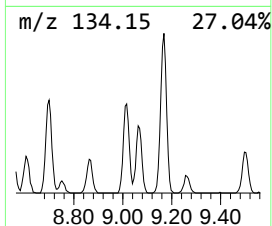
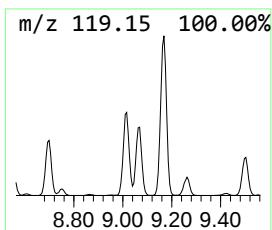
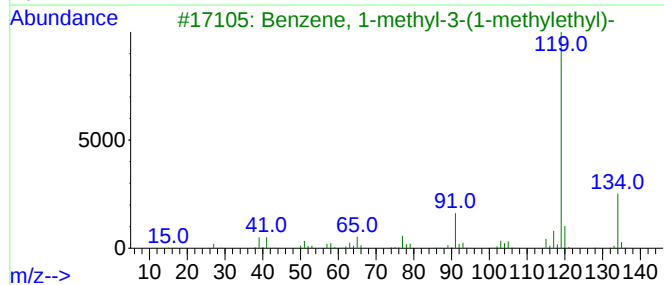
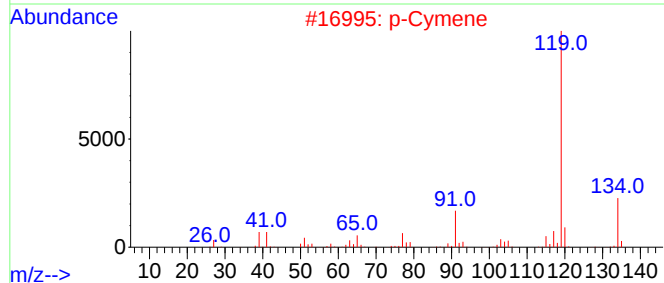
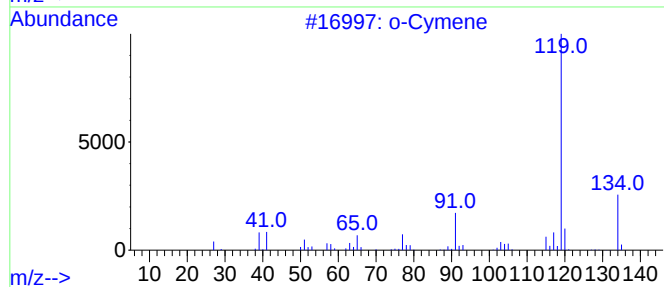
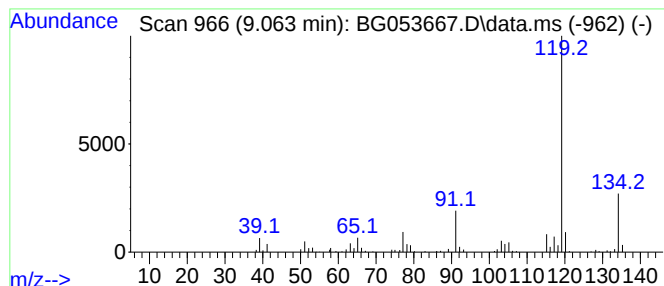
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 o-Cymene Concentration Rank 18

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 9.063 | 26.52 ng | 273032 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 97   |
| 2    |    |   | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 97   |
| 3    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 95   |
| 4    |    |   | Benzene, 1,2,3,5-tetramethyl-       | 134 | C10H14  | 000527-53-7 | 94   |
| 5    |    |   | 1,3,5-Cycloheptatriene, 3,7,7-tr... | 134 | C10H14  | 003479-89-8 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

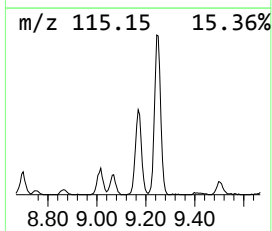
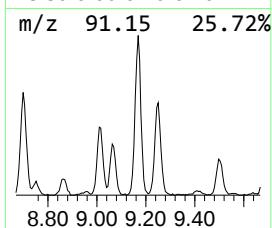
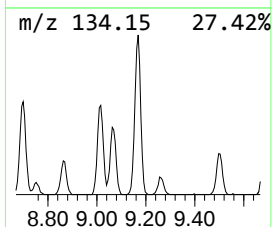
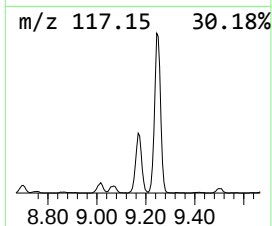
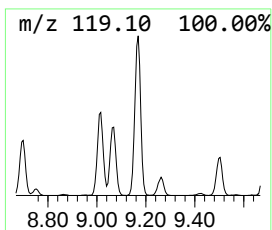
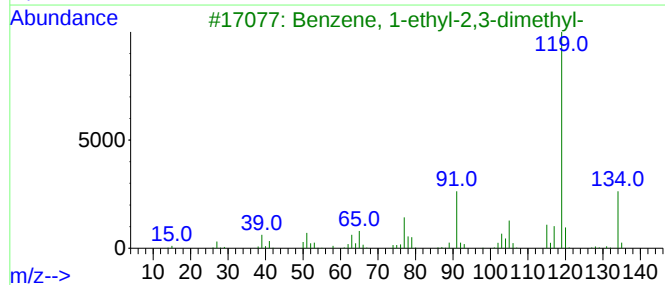
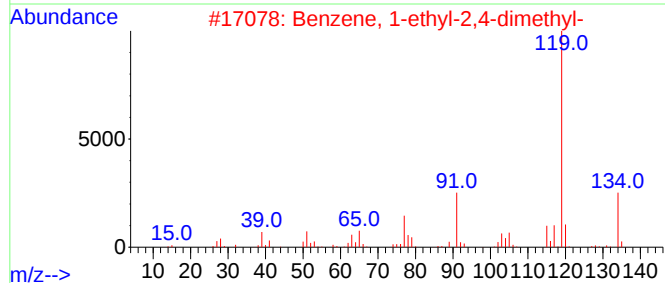
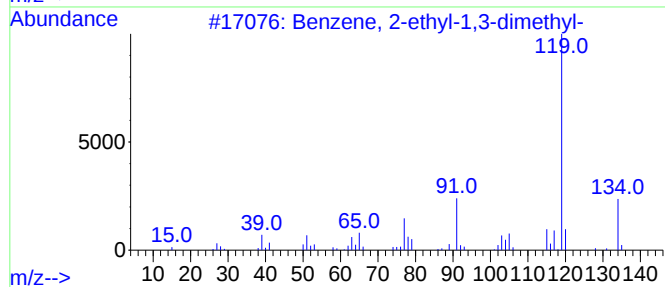
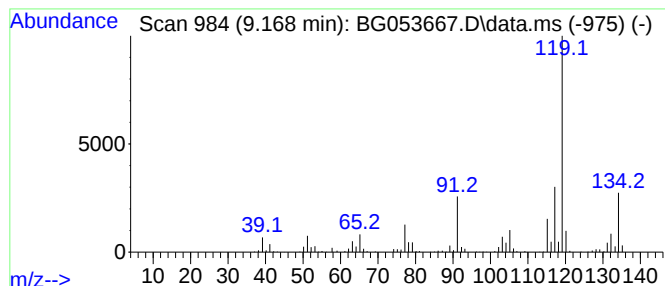
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 10

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 9.168 | 86.29 ng | 888285 | 1,4-Dichlorobenzene-d4 | 8.058 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 94   |
| 2    |    |   | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 94   |
| 3    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 93   |
| 4    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 93   |
| 5    |    |   | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

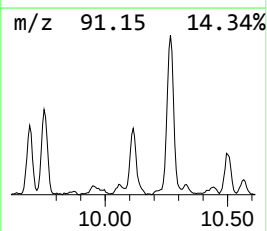
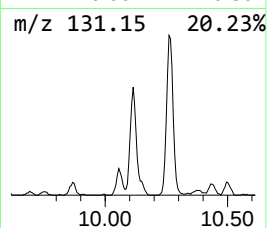
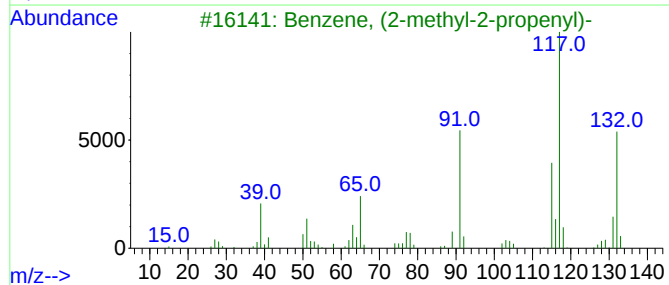
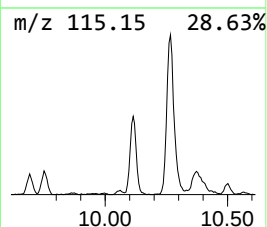
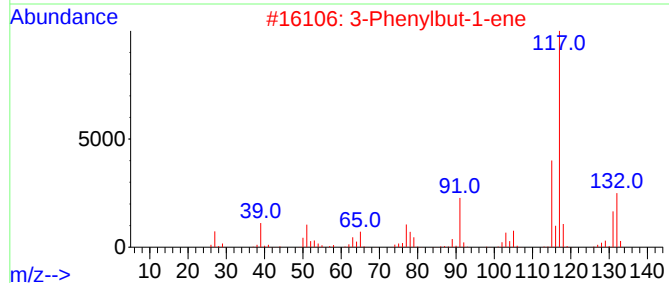
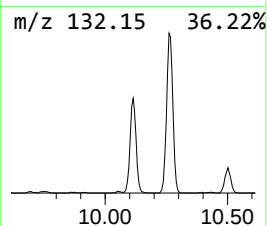
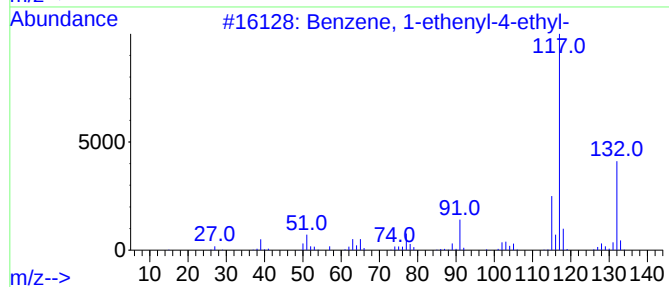
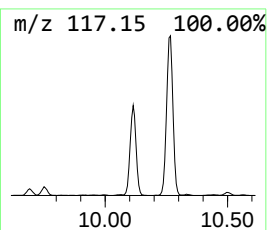
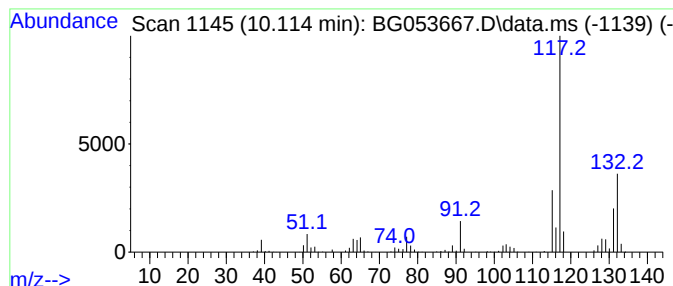
Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 17

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 10.114    | 26.91 ng                            | 701836 | Naphthalene-d8   | 10.872      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Benzene, 1-ethenyl-4-ethyl-         | 132    | C10H12           | 003454-07-7 | 94   |
| 2         | 3-Phenylbut-1-ene                   | 132    | C10H12           | 000934-10-1 | 91   |
| 3         | Benzene, (2-methyl-2-propenyl)-     | 132    | C10H12           | 003290-53-7 | 90   |
| 4         | Benzene, 1-ethenyl-3-ethyl-         | 132    | C10H12           | 007525-62-4 | 87   |
| 5         | Benzene, (1-methyl-1-propenyl)-,... | 132    | C10H12           | 000768-00-3 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

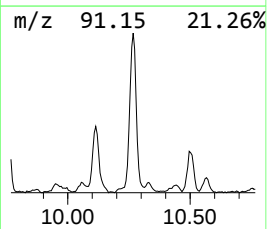
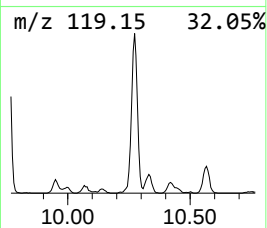
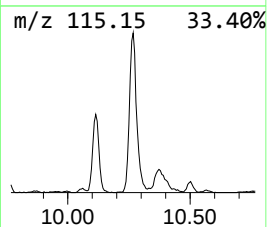
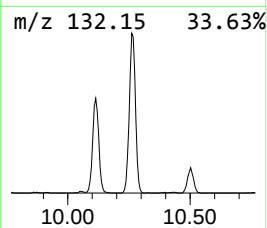
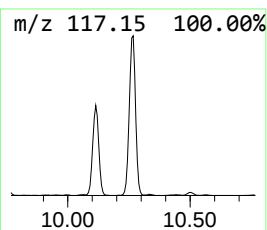
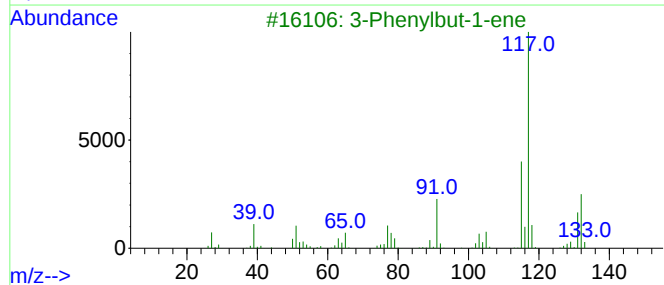
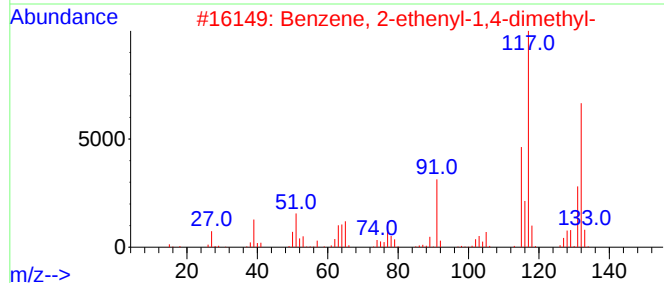
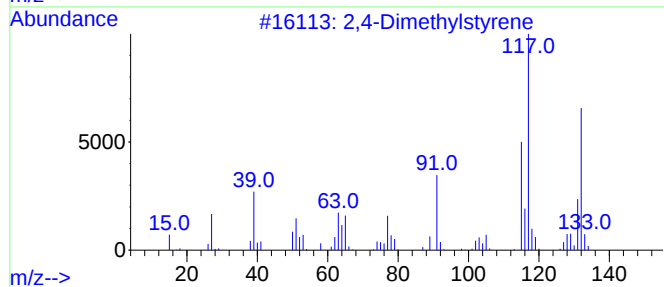
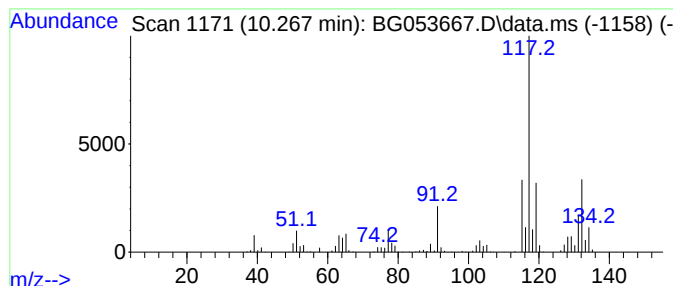
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 2,4-Dimethylstyrene Concentration Rank 12

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 10.267 | 65.15 ng | 1699370 | Naphthalene-d8   | 10.872 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2,4-Dimethylstyrene               | 132 | C10H12  | 002234-20-0 | 95   |
| 2    |    |   | Benzene, 2-ethenyl-1,4-dimethyl-  | 132 | C10H12  | 002039-89-6 | 95   |
| 3    |    |   | 3-Phenylbut-1-ene                 | 132 | C10H12  | 000934-10-1 | 81   |
| 4    |    |   | Benzene, 1-methyl-4-(2-propenyl)- | 132 | C10H12  | 003333-13-9 | 78   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-4-methyl-  | 132 | C10H12  | 000824-22-6 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG052322\  
Data File : BG053667.D  
Acq On : 23 May 2022 16:49  
Operator : CG/JU  
Sample : N2968-04  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-25

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |        |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|--------|------|
|                    |        |         |       |          | #                     | RT     | Resp   | Conc |
| Di-sec-Butyl ether | 4.533  | 22.7    | ng    | 234037   | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, (1-met... | 6.543  | 27.5    | ng    | 283228   | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, propyl-   | 7.030  | 66.0    | ng    | 679112   | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 1-ethy... | 7.142  | 139.7   | ng    | 1437870  | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 1-ethy... | 7.447  | 165.7   | ng    | 1705700  | 1                     | 8.058  | 205879 | 20.0 |
| Mesitylene         | 7.712  | 504.1   | ng    | 5188780  | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 1,2,3-... | 8.176  | 176.1   | ng    | 1813020  | 1                     | 8.058  | 205879 | 20.0 |
| Indane             | 8.434  | 218.2   | ng    | 2245630  | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 1,2-di... | 8.546  | 22.4    | ng    | 230850   | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 1-prop... | 8.604  | 32.6    | ng    | 335716   | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 1,4-di... | 8.698  | 39.2    | ng    | 403755   | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 2-ethy... | 9.016  | 35.8    | ng    | 368399   | 1                     | 8.058  | 205879 | 20.0 |
| o-Cymene           | 9.063  | 26.5    | ng    | 273032   | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 2-ethy... | 9.168  | 86.3    | ng    | 888285   | 1                     | 8.058  | 205879 | 20.0 |
| Benzene, 1-ethe... | 10.114 | 26.9    | ng    | 701836   | 2                     | 10.872 | 521644 | 20.0 |
| 2,4-Dimethylsty... | 10.267 | 65.2    | ng    | 1699370  | 2                     | 10.872 | 521644 | 20.0 |