

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
 Data File : BG057840.D  
 Acq On : 23 Jun 2023 13:22  
 Operator : CG/JU  
 Sample : 03291-03  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MW-6A

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-BG061923.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG057840.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         | 4.128       | 167           | 177         | 179          | rBV2     | 9554199        | 22961819      | 100.00%         | 26.510%       |
| 2         | 5.433       | 391           | 399         | 406          | rBV      | 2119514        | 3419226       | 14.89%          | 3.948%        |
| 3         | 5.503       | 406           | 411         | 415          | rVV      | 378909         | 566328        | 2.47%           | 0.654%        |
| 4         | 5.580       | 415           | 424         | 433          | rVB      | 6371744        | 17081028      | 74.39%          | 19.721%       |
| 5         | 5.938       | 477           | 485         | 495          | rVB      | 4374804        | 7747501       | 33.74%          | 8.945%        |
| 6         | 7.013       | 658           | 668         | 673          | rVV      | 1696999        | 2805993       | 12.22%          | 3.240%        |
| 7         | 7.072       | 673           | 678         | 685          | rVV2     | 945378         | 1647928       | 7.18%           | 1.903%        |
| 8         | 7.143       | 685           | 690         | 700          | rVB      | 838267         | 1361833       | 5.93%           | 1.572%        |
| 9         | 7.313       | 710           | 719         | 725          | rBV      | 866077         | 1400406       | 6.10%           | 1.617%        |
| 10        | 7.577       | 752           | 764         | 773          | rBV      | 2694707        | 4728188       | 20.59%          | 5.459%        |
| 11        | 7.924       | 815           | 823         | 827          | rBV      | 141607         | 253083        | 1.10%           | 0.292%        |
| 12        | 8.036       | 835           | 842         | 849          | rVB      | 970918         | 1574657       | 6.86%           | 1.818%        |
| 13        | 8.300       | 879           | 887         | 892          | rBV      | 956104         | 1626937       | 7.09%           | 1.878%        |
| 14        | 8.365       | 892           | 898         | 903          | rVV      | 685356         | 1103678       | 4.81%           | 1.274%        |
| 15        | 8.465       | 910           | 915         | 922          | rVB2     | 231069         | 410524        | 1.79%           | 0.474%        |
| 16        | 8.559       | 924           | 931         | 936          | rBV      | 267796         | 447744        | 1.95%           | 0.517%        |
| 17        | 9.029       | 1003          | 1011        | 1015         | rBV      | 291237         | 487014        | 2.12%           | 0.562%        |
| 18        | 9.087       | 1015          | 1021        | 1033         | rVB2     | 481527         | 1172667       | 5.11%           | 1.354%        |
| 19        | 9.552       | 1094          | 1100        | 1105         | rBV      | 153075         | 242424        | 1.06%           | 0.280%        |
| 20        | 9.610       | 1105          | 1110        | 1116         | rVB      | 196757         | 322645        | 1.41%           | 0.373%        |
| 21        | 9.687       | 1116          | 1123        | 1132         | rBV      | 267171         | 426435        | 1.86%           | 0.492%        |
| 22        | 9.969       | 1166          | 1171        | 1181         | rVB      | 220612         | 405840        | 1.77%           | 0.469%        |
| 23        | 10.121      | 1191          | 1197        | 1207         | rBV2     | 382690         | 752376        | 3.28%           | 0.869%        |
| 24        | 10.727      | 1289          | 1300        | 1303         | rBV2     | 189902         | 408130        | 1.78%           | 0.471%        |
| 25        | 10.779      | 1303          | 1309        | 1316         | rVB      | 1371774        | 2469618       | 10.76%          | 2.851%        |
| 26        | 10.897      | 1323          | 1329        | 1334         | rBV      | 163447         | 274891        | 1.20%           | 0.317%        |
| 27        | 12.384      | 1576          | 1582        | 1588         | rBV      | 229685         | 362493        | 1.58%           | 0.419%        |
| 28        | 12.601      | 1610          | 1619        | 1629         | rVB2     | 150334         | 304526        | 1.33%           | 0.352%        |
| 29        | 13.183      | 1709          | 1718        | 1725         | rBV      | 1223243        | 1869212       | 8.14%           | 2.158%        |
| 30        | 14.557      | 1943          | 1952        | 1960         | rBV      | 351705         | 543587        | 2.37%           | 0.628%        |
| 31        | 15.915      | 2178          | 2183        | 2190         | rVB      | 267171         | 383565        | 1.67%           | 0.443%        |
| 32        | 16.044      | 2197          | 2205        | 2216         | rBV      | 1200038        | 1849544       | 8.05%           | 2.135%        |
| 33        | 17.301      | 2413          | 2419        | 2425         | rBV      | 474145         | 707101        | 3.08%           | 0.816%        |
| 34        | 19.916      | 2858          | 2864        | 2870         | rBV      | 2061775        | 2918193       | 12.71%          | 3.369%        |
| 35        | 21.555      | 3137          | 3143        | 3151         | rVB2     | 453739         | 731140        | 3.18%           | 0.844%        |
| 36        | 24.710      | 3669          | 3680        | 3691         | rBV      | 298161         | 847082        | 3.69%           | 0.978%        |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

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-BG061923.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

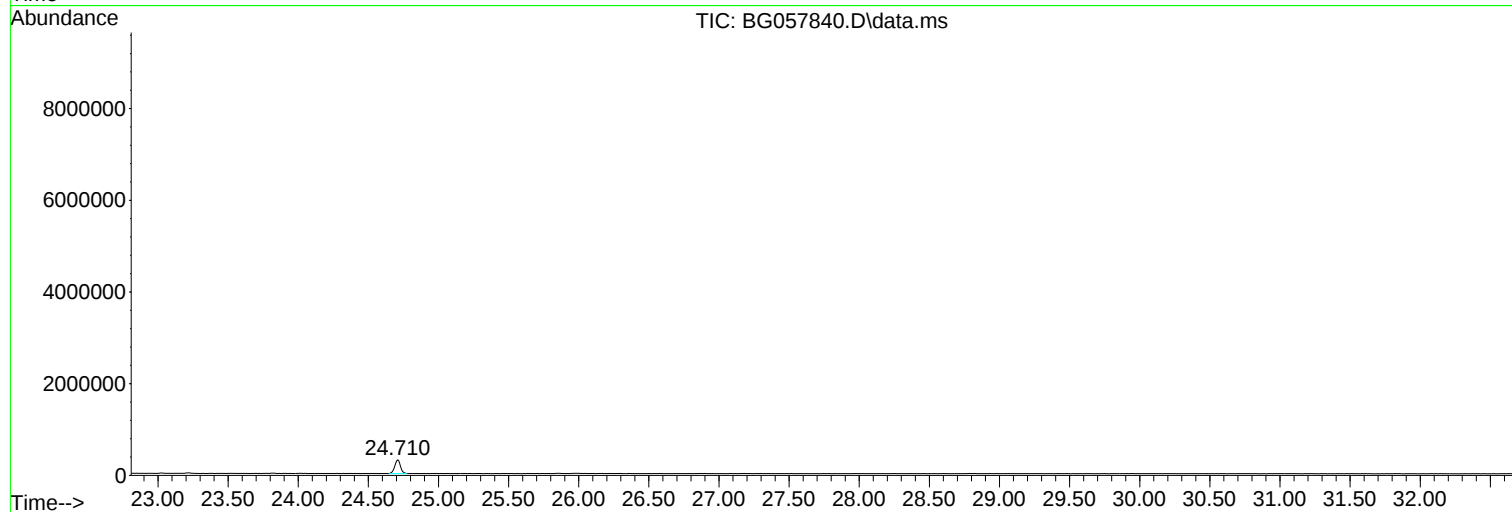
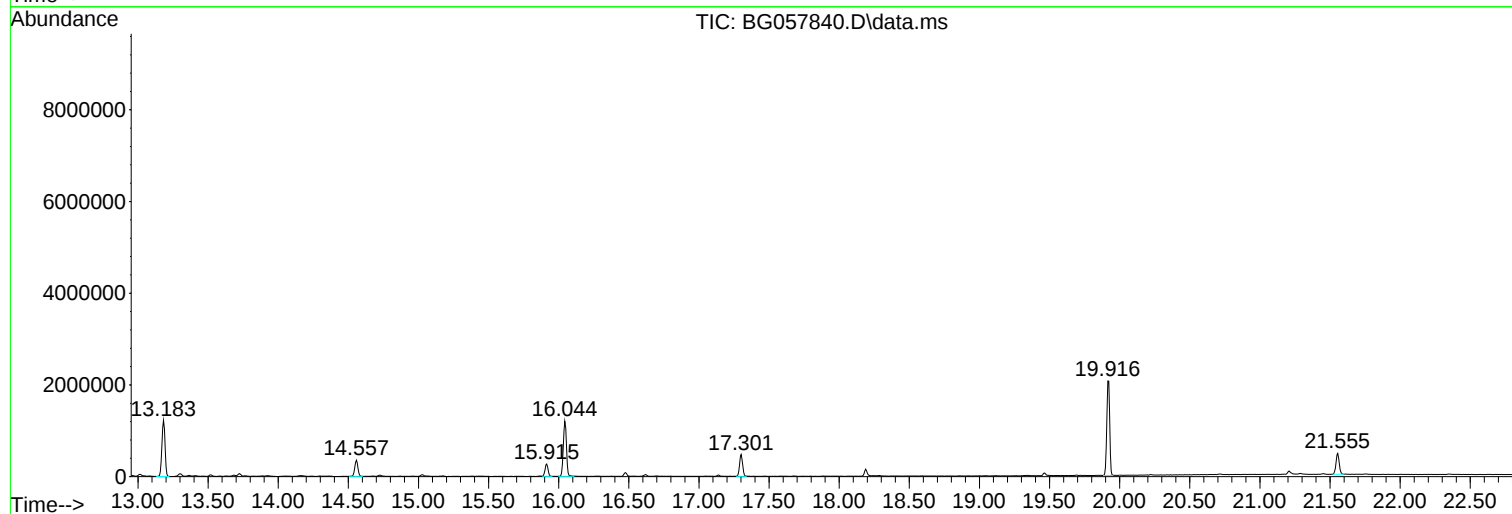
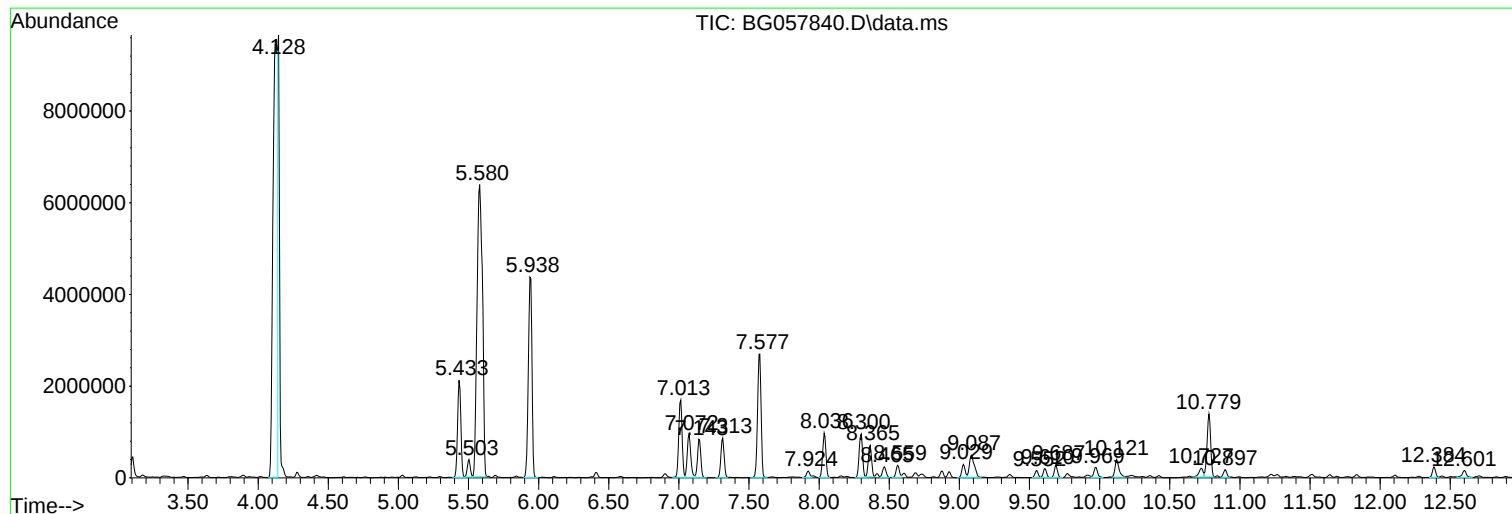
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.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\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

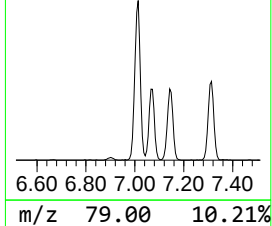
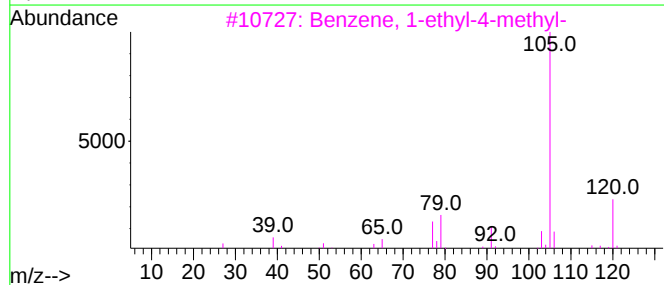
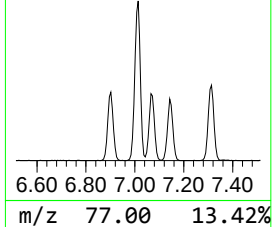
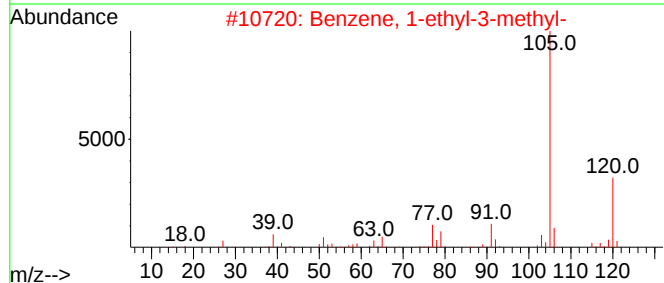
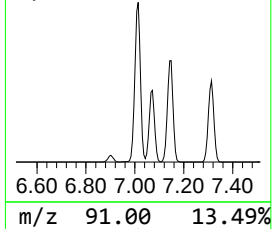
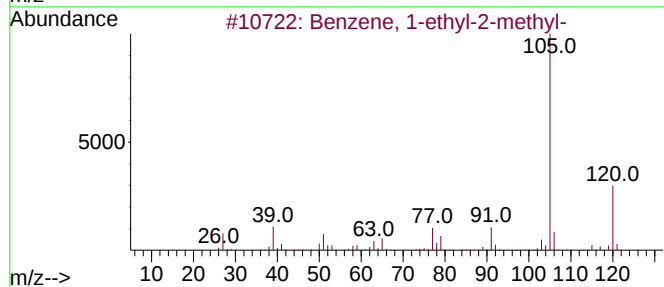
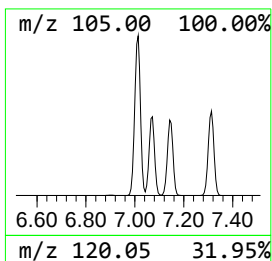
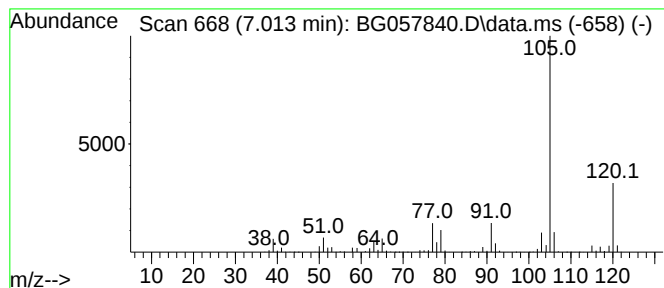
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 7.013 | 221.75 ng | 2805990 | 1,4-Dichlorobenzene-d4 | 7.924 |

| 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 | 95   |
| 3    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 93   |
| 4    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

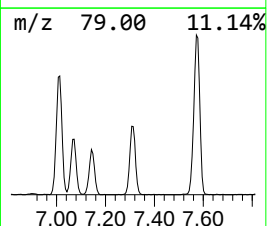
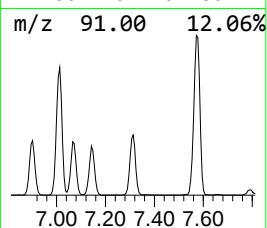
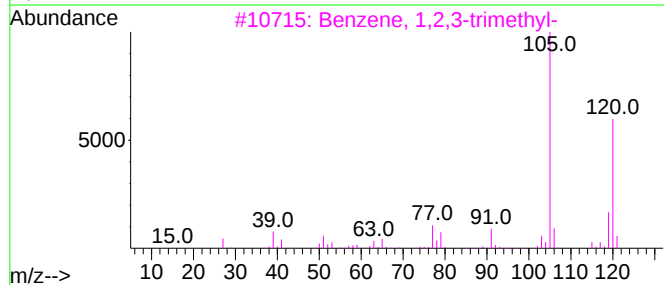
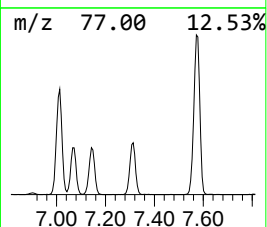
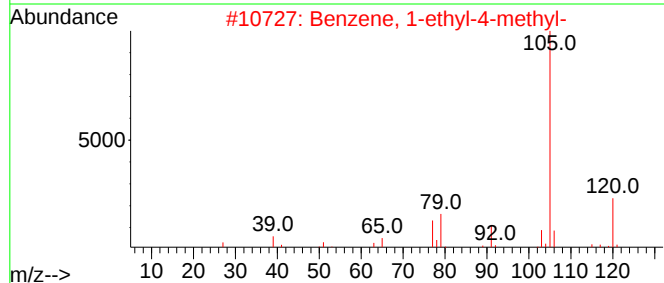
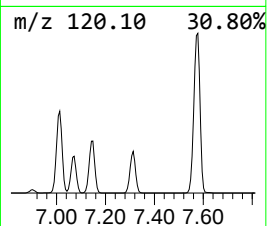
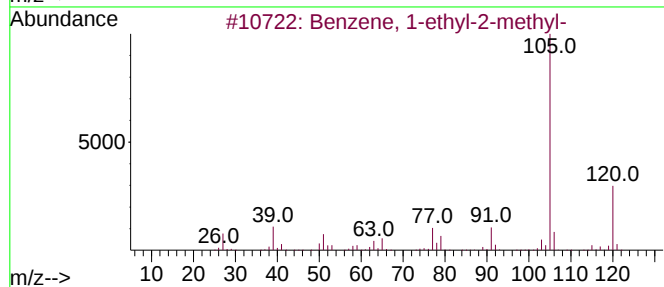
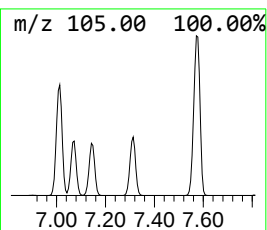
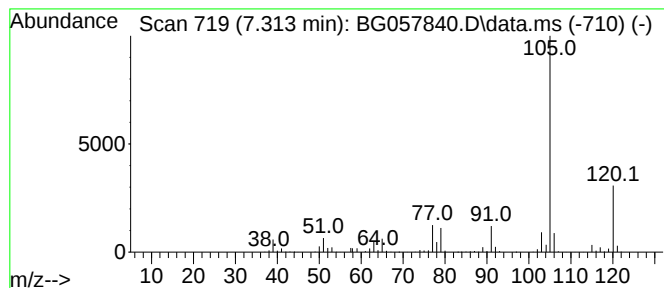
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Benzene, 1-ethyl-4-methyl- Concentration Rank 9

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 7.313 | 110.67 ng | 1400410 | 1,4-Dichlorobenzene-d4 | 7.924 |

| 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 | 94   |
| 3    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

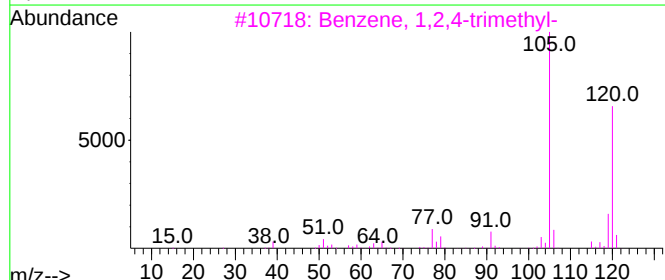
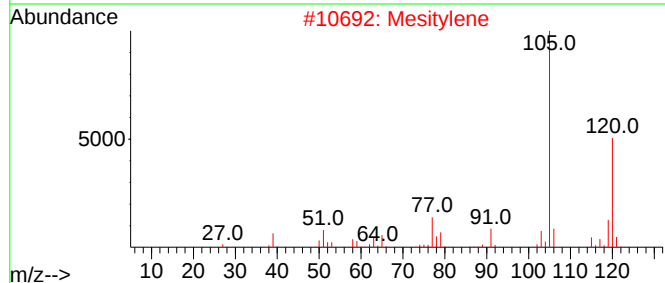
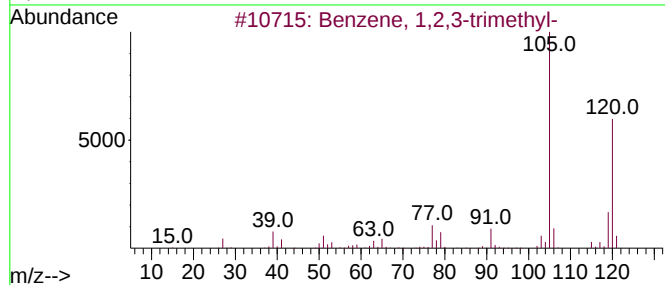
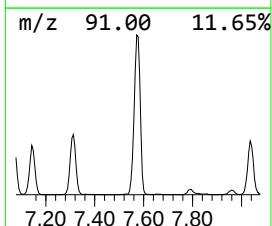
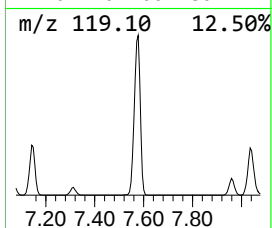
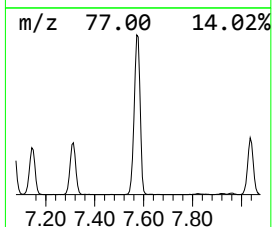
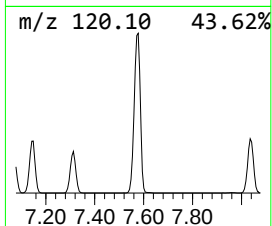
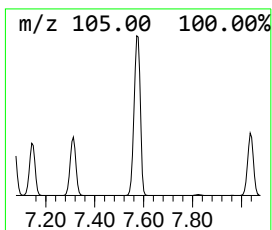
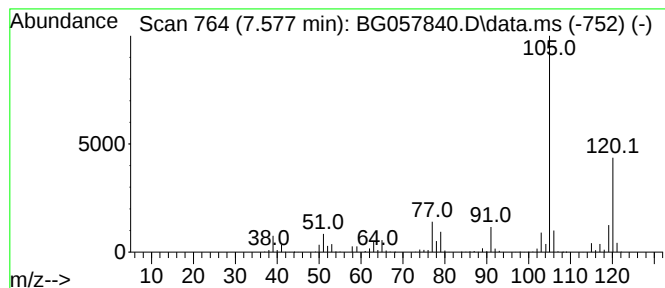
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 4

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 7.577 | 373.65 ng | 4728190 | 1,4-Dichlorobenzene-d4 | 7.924 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 97   |
| 3    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
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Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

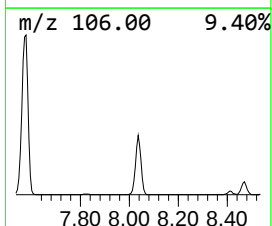
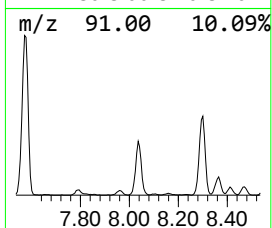
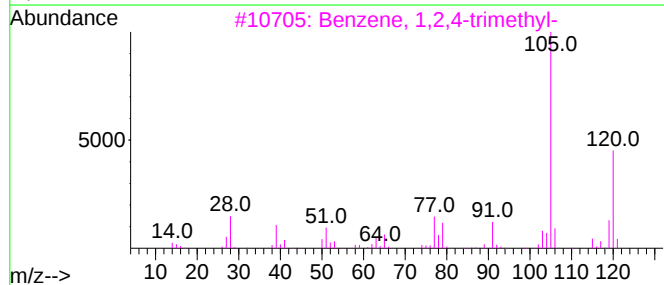
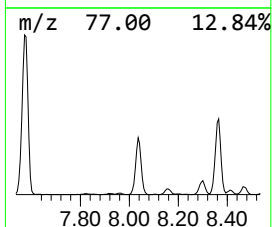
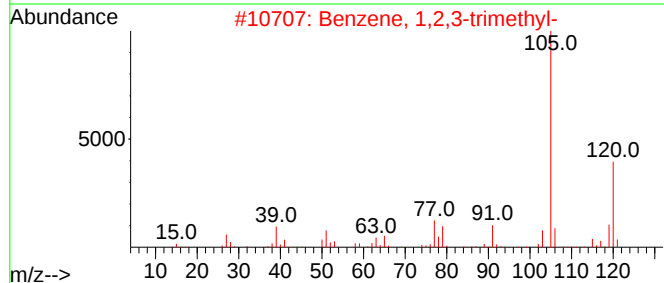
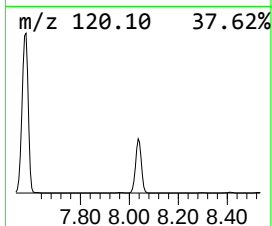
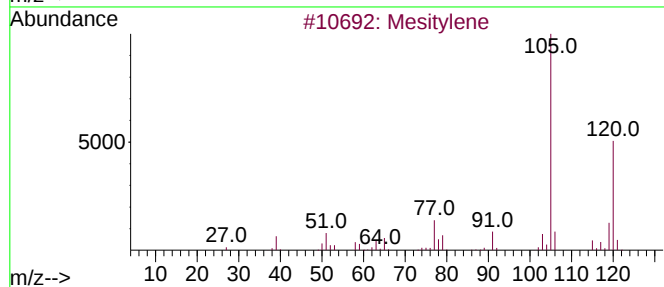
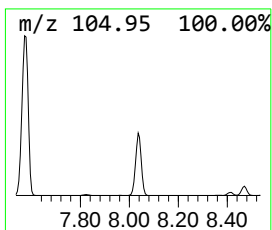
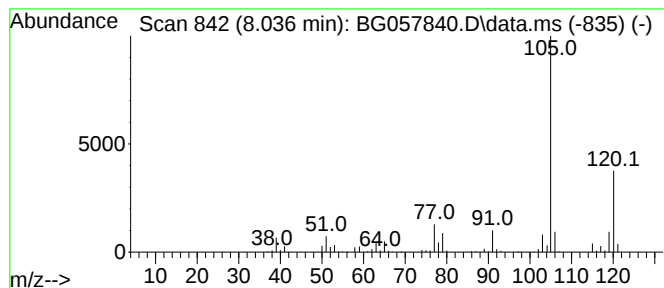
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Mesitylene Concentration Rank 8

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 8.036 | 124.44 ng | 1574660 | 1,4-Dichlorobenzene-d4 | 7.924 |

| 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 | 95   |
| 3    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 90   |
| 5    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 87   |



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Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

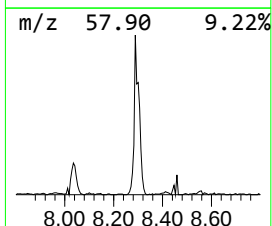
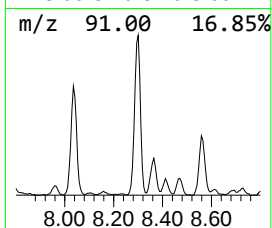
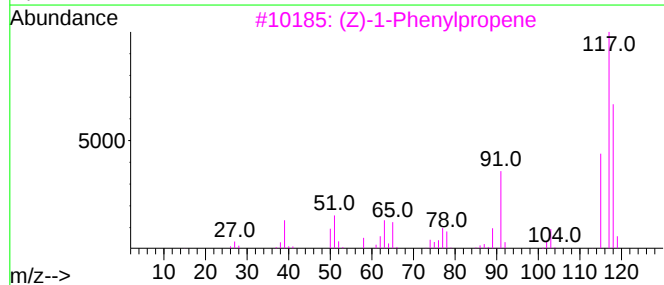
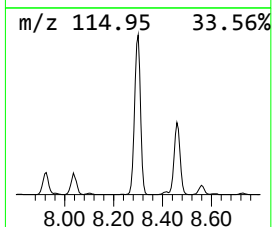
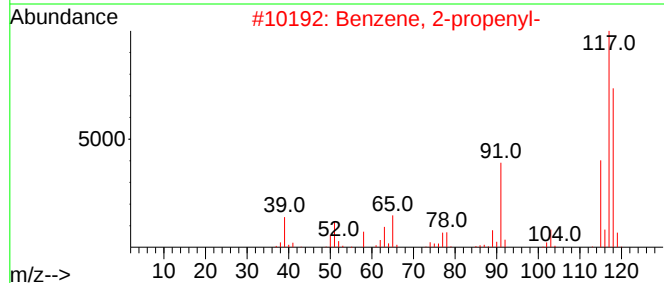
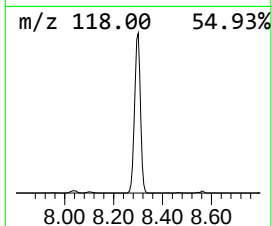
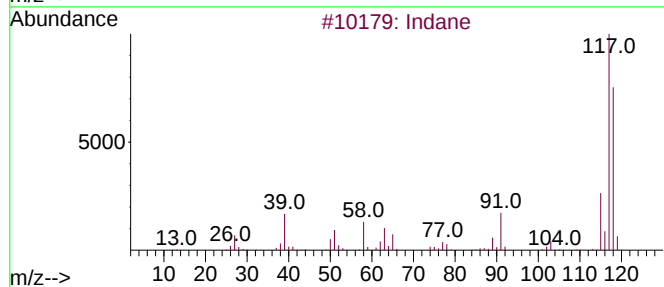
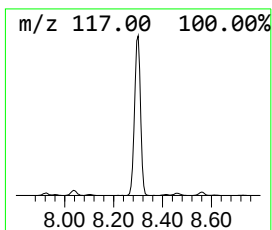
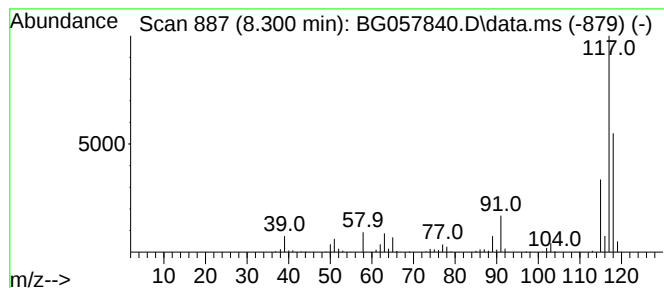
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Indane Concentration Rank 7

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 8.300 | 128.57 ng | 1626940 | 1,4-Dichlorobenzene-d4 | 7.924 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Indane                              | 118 | C9H10   | 000496-11-7  | 92   |
| 2    |    |   | Benzene, 2-propenyl-                | 118 | C9H10   | 000300-57-2  | 81   |
| 3    |    |   | (Z)-1-Phenylpropene                 | 118 | C9H10   | 000766-90-5  | 76   |
| 4    |    |   | Delatacylene                        | 118 | C9H10   | 007785-10-6  | 74   |
| 5    |    |   | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10   | 1000191-13-7 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

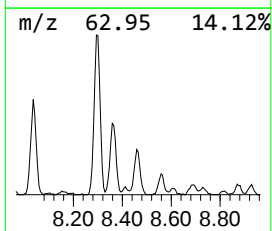
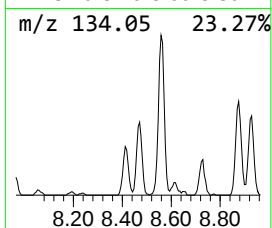
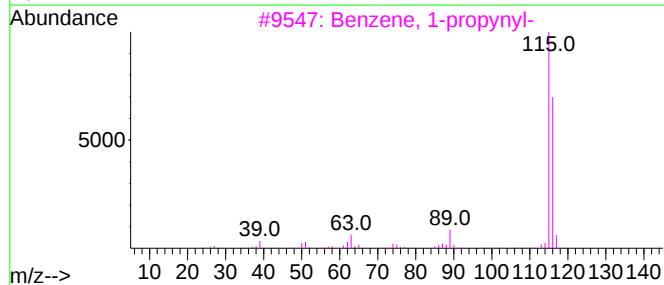
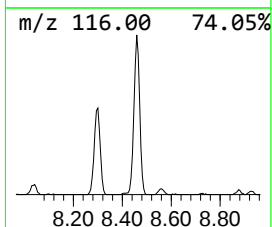
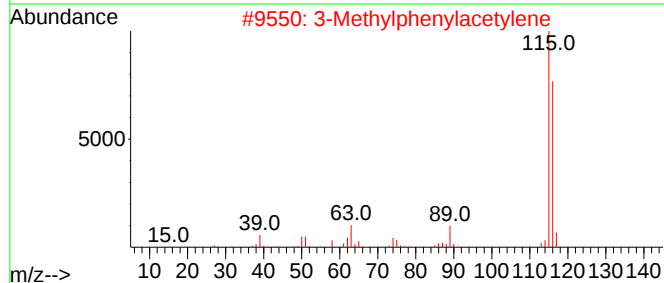
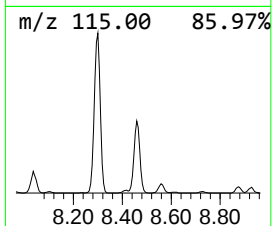
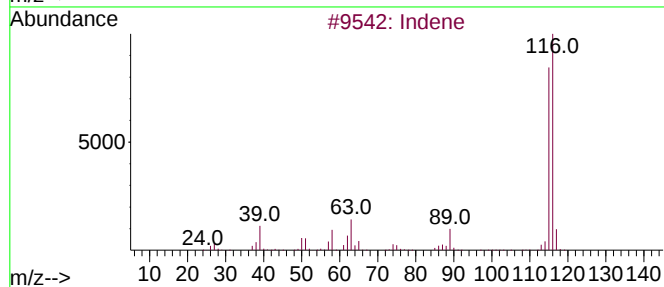
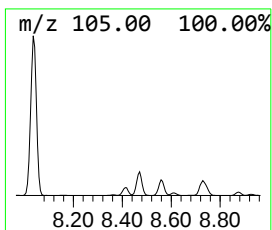
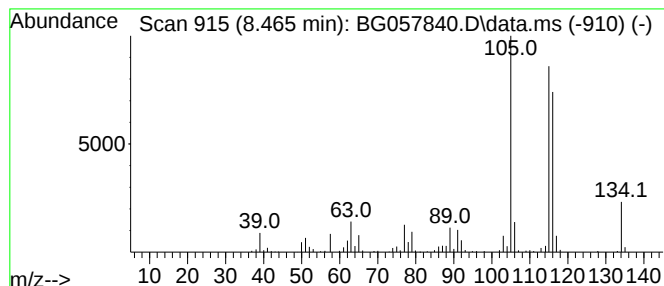
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 Indene Concentration Rank 13

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 8.465 | 32.44 ng | 410524 | 1,4-Dichlorobenzene-d4 | 7.924 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------|-----|---------|-------------|------|
| 1    |      | Indene                       | 116 | C9H8    | 000095-13-6 | 93   |
| 2    |      | 3-Methylphenylacetylene      | 116 | C9H8    | 000766-82-5 | 76   |
| 3    |      | Benzene, 1-propynyl-         | 116 | C9H8    | 000673-32-5 | 76   |
| 4    |      | 2-Methylphenylacetylene      | 116 | C9H8    | 000766-47-2 | 64   |
| 5    |      | Benzene, 1-ethynyl-4-methyl- | 116 | C9H8    | 000766-97-2 | 62   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

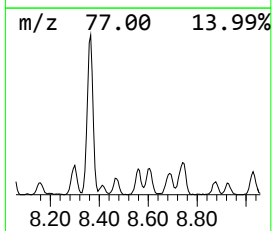
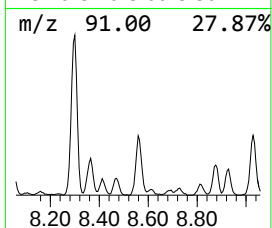
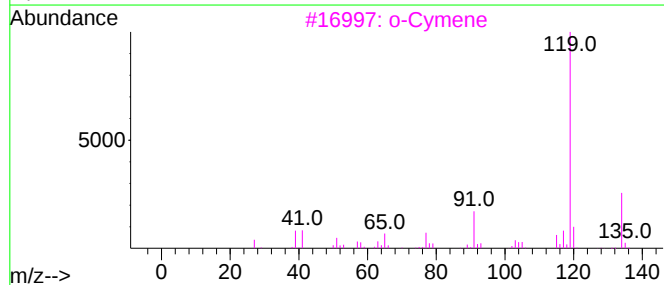
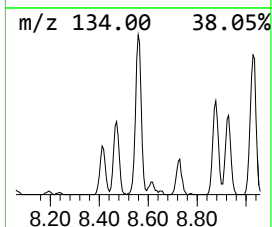
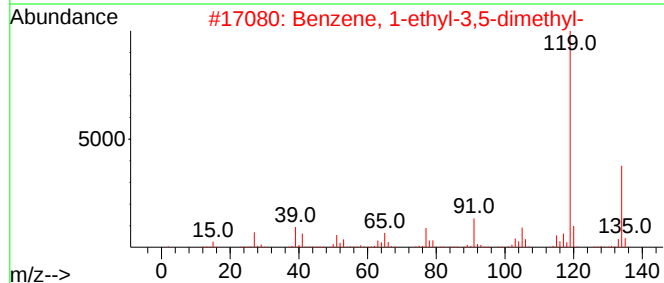
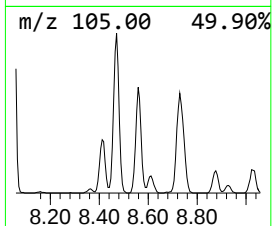
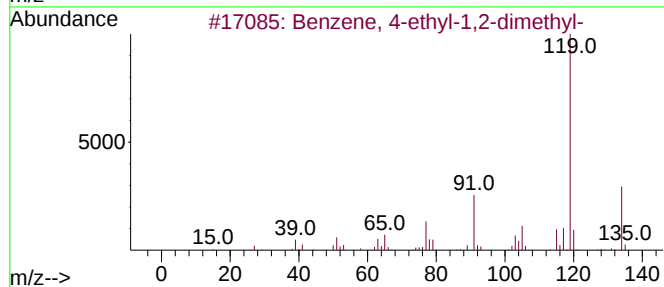
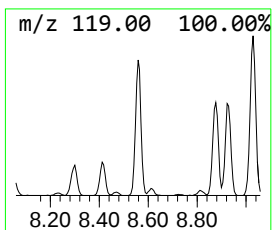
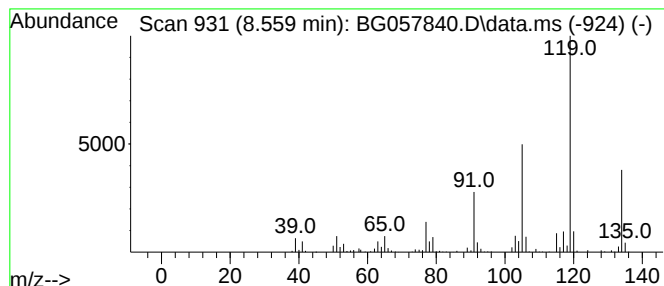
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 12

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 8.559 | 35.38 ng | 447744 | 1,4-Dichlorobenzene-d4 | 7.924 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 95   |
| 2    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 95   |
| 3    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 93   |
| 4    |    |   | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 93   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

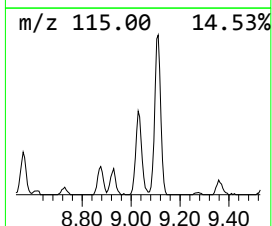
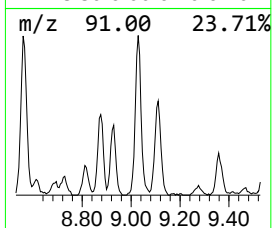
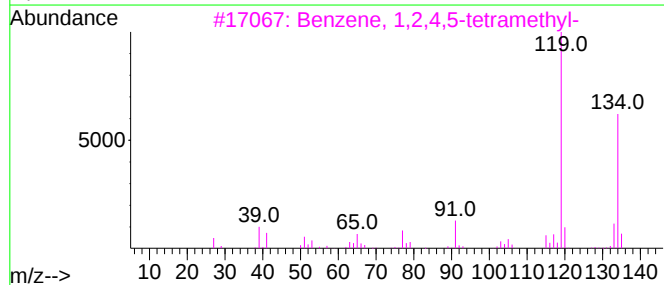
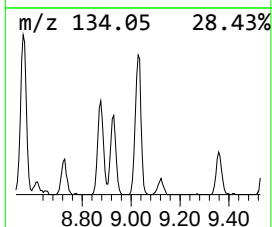
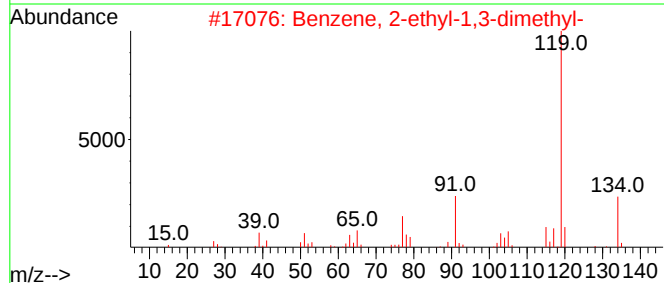
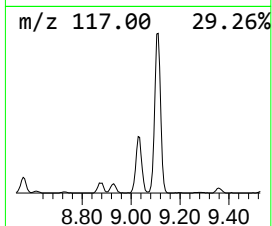
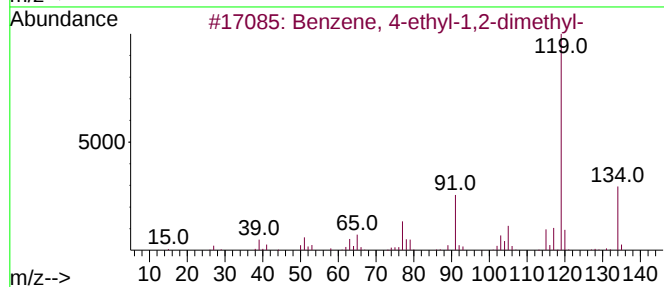
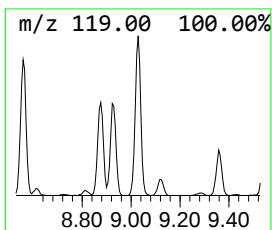
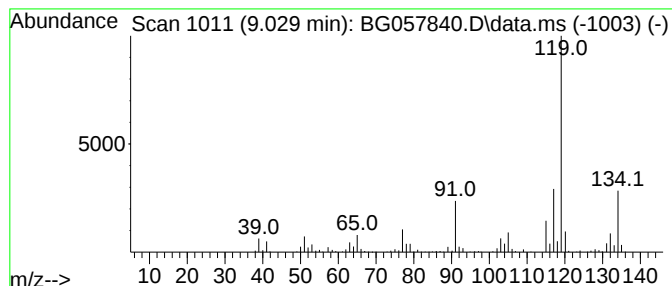
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 10

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 9.029 | 38.49 ng | 487014 | 1,4-Dichlorobenzene-d4 | 7.924 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 94   |
| 2    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 90   |
| 3    |    |   | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 83   |
| 4    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 81   |
| 5    |    |   | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

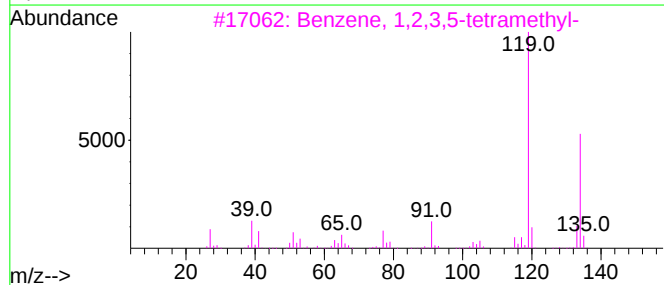
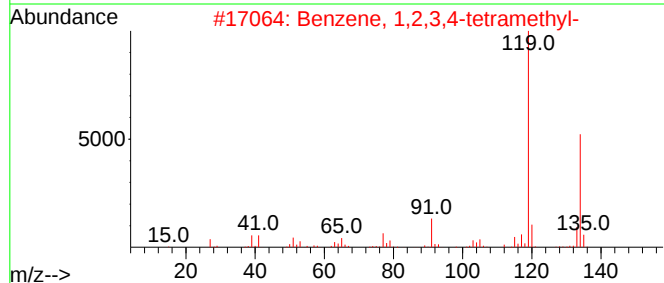
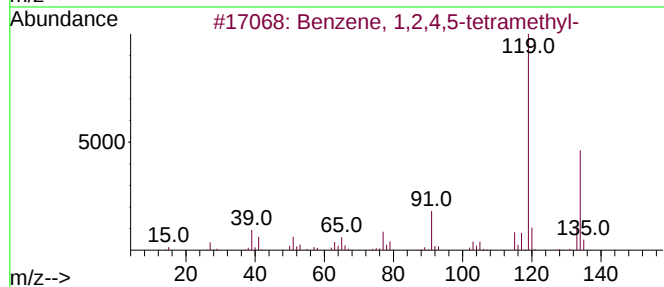
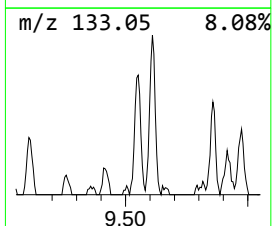
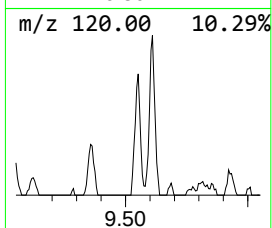
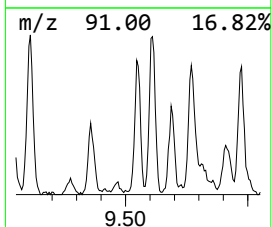
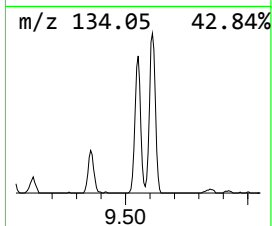
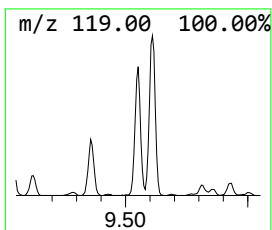
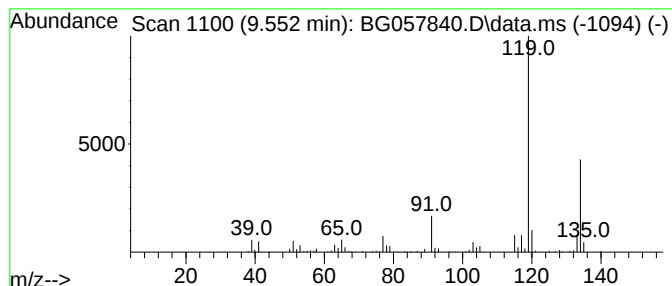
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 19

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.   |
|-------|----------|--------|------------------|--------|
| 9.552 | 11.88 ng | 242424 | Naphthalene-d8   | 10.727 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-       | 134 | C10H14  | 000095-93-2 | 97   |
| 2    |    |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 95   |
| 3    |    |   | Benzene, 1,2,3,5-tetramethyl-       | 134 | C10H14  | 000527-53-7 | 95   |
| 4    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 95   |
| 5    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

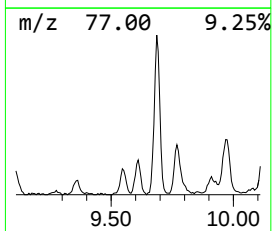
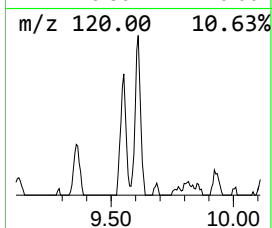
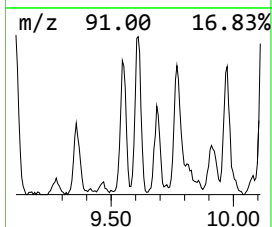
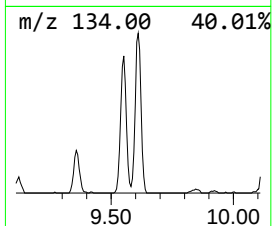
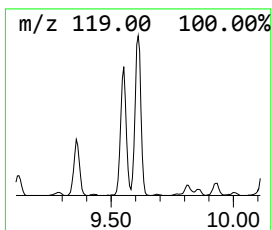
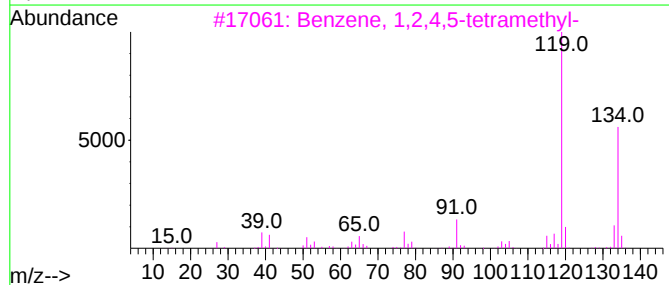
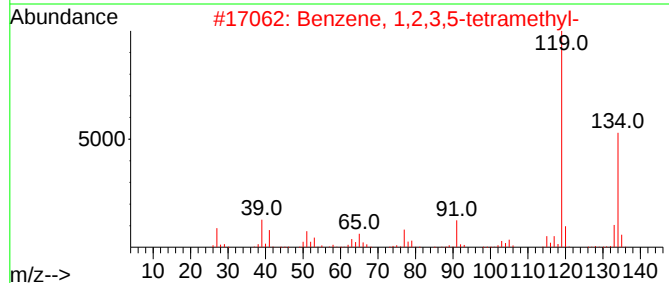
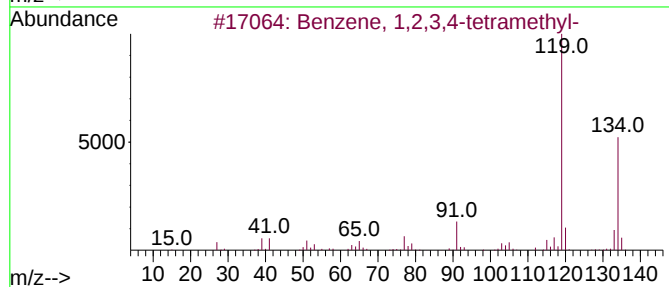
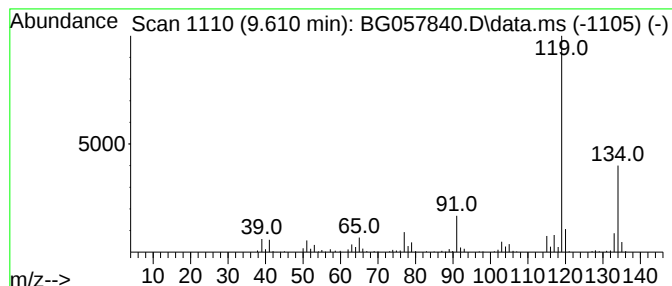
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 16

| R.T.      | EstConc                        | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------|--------|------------------|-------------|------|
| 9.610     | 15.81 ng                       | 322645 | Naphthalene-d8   | 10.727      |      |
| Hit# of 5 | Tentative ID                   | MW     | MolForm          | CAS#        | Qual |
| 1         | Benzene, 1,2,3,4-tetramethyl-  | 134    | C10H14           | 000488-23-3 | 95   |
| 2         | Benzene, 1,2,3,5-tetramethyl-  | 134    | C10H14           | 000527-53-7 | 95   |
| 3         | Benzene, 1,2,4,5-tetramethyl-  | 134    | C10H14           | 000095-93-2 | 95   |
| 4         | o-Cymene                       | 134    | C10H14           | 000527-84-4 | 95   |
| 5         | Benzene, 1-ethyl-2,3-dimethyl- | 134    | C10H14           | 000933-98-2 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

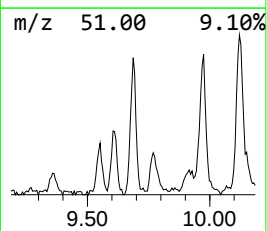
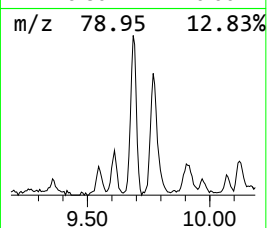
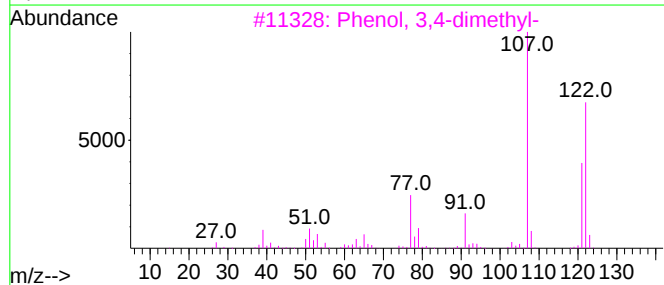
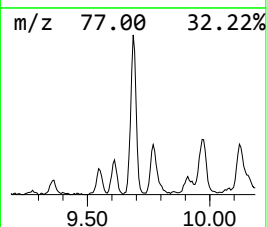
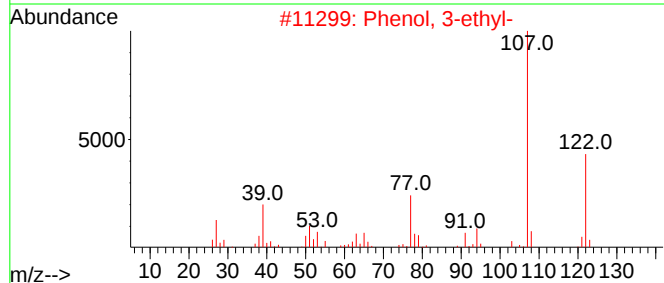
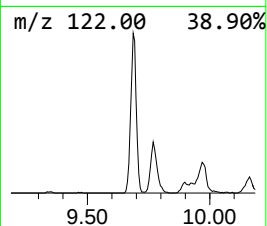
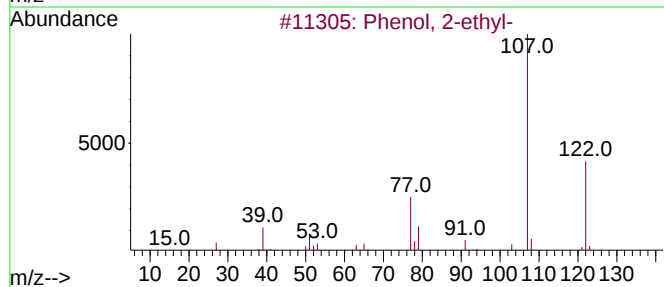
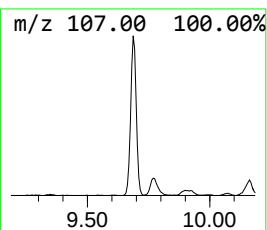
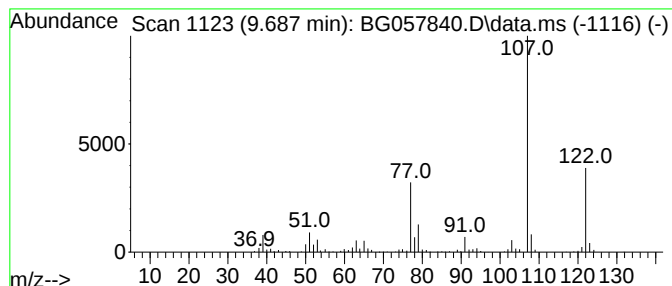
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 Phenol, 2-ethyl- Concentration Rank 14

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.   |
|-------|----------|--------|------------------|--------|
| 9.687 | 20.90 ng | 426435 | Naphthalene-d8   | 10.727 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Phenol, 2-ethyl-      | 122 | C8H10O  | 000090-00-6 | 95   |
| 2    |    |   | Phenol, 3-ethyl-      | 122 | C8H10O  | 000620-17-7 | 91   |
| 3    |    |   | Phenol, 3,4-dimethyl- | 122 | C8H10O  | 000095-65-8 | 90   |
| 4    |    |   | Phenol, 4-ethyl-      | 122 | C8H10O  | 000123-07-9 | 87   |
| 5    |    |   | 2-Isopropylpyrazine   | 122 | C7H10N2 | 029460-90-0 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

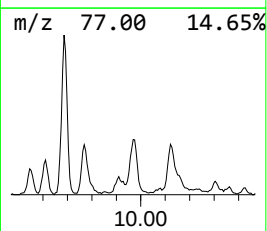
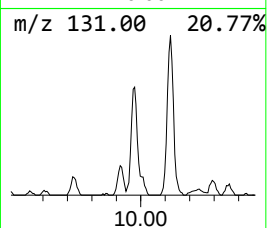
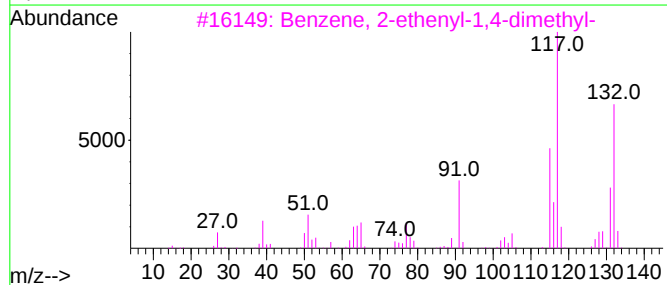
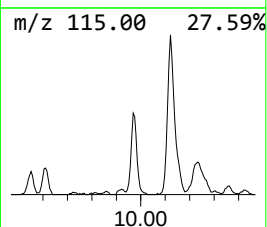
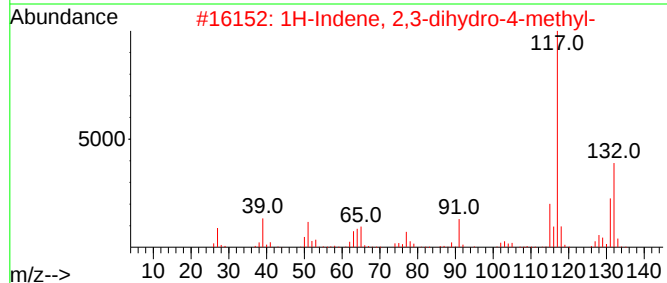
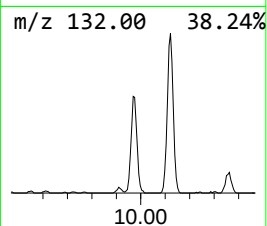
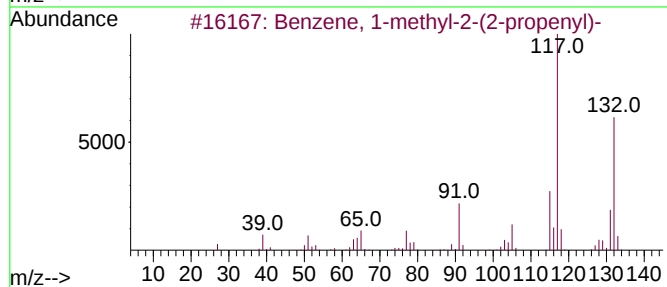
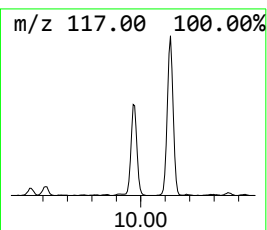
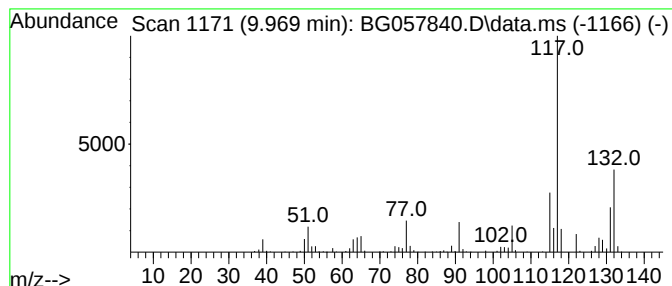
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, 1-methyl-2-(2-prop... Concentration Rank 15

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.   |
|-------|----------|--------|------------------|--------|
| 9.969 | 19.89 ng | 405840 | Naphthalene-d8   | 10.727 |

| Hit# | of 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------------------|-----|---------|-------------|------|
| 1    |      | Benzene, 1-methyl-2-(2-propenyl)- | 132 | C10H12  | 001587-04-8 | 90   |
| 2    |      | 1H-Indene, 2,3-dihydro-4-methyl-  | 132 | C10H12  | 000824-22-6 | 87   |
| 3    |      | Benzene, 2-ethenyl-1,4-dimethyl-  | 132 | C10H12  | 002039-89-6 | 86   |
| 4    |      | 3-Phenylbut-1-ene                 | 132 | C10H12  | 000934-10-1 | 86   |
| 5    |      | Indan, 1-methyl-                  | 132 | C10H12  | 000767-58-8 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

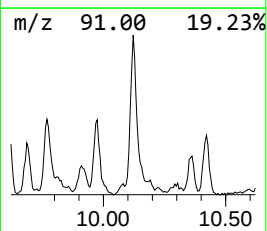
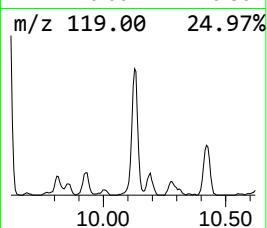
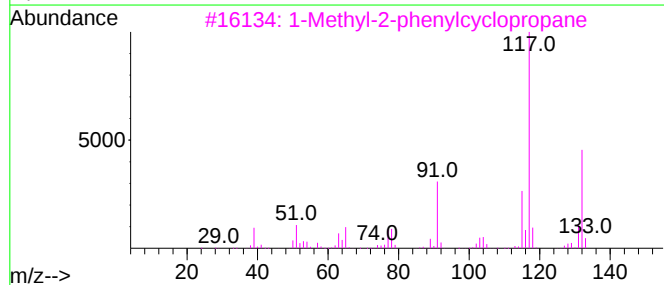
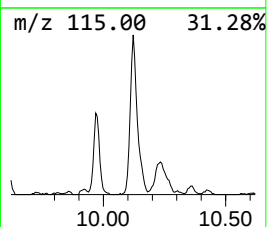
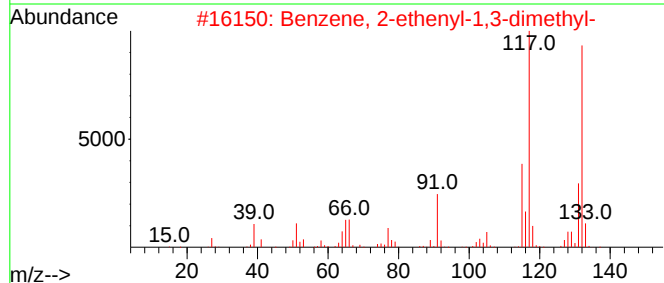
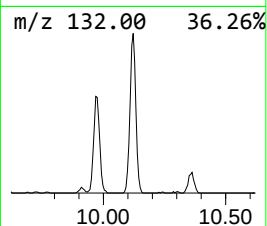
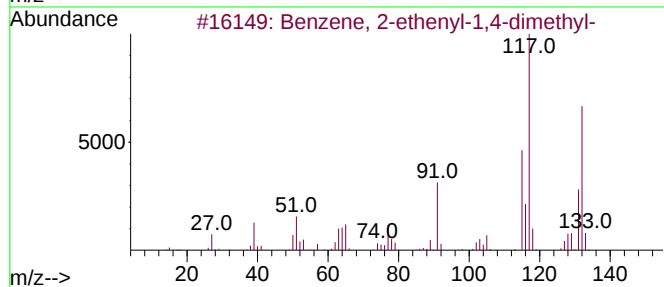
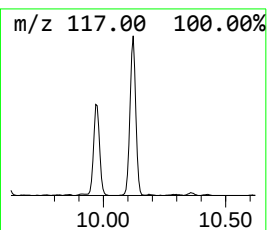
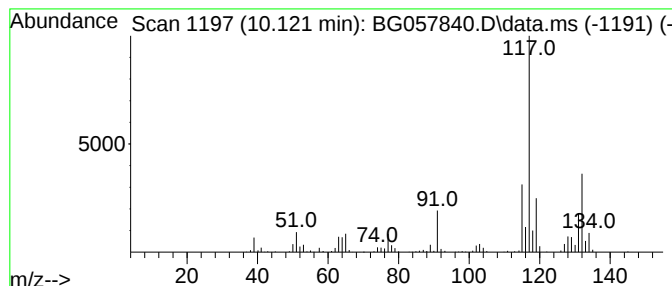
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 10.121 | 36.87 ng | 752376 | Naphthalene-d8   | 10.727 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12  | 002039-89-6 | 96   |
| 2    |    |   | Benzene, 2-ethenyl-1,3-dimethyl- | 132 | C10H12  | 002039-90-9 | 93   |
| 3    |    |   | 1-Methyl-2-phenylcyclopropane    | 132 | C10H12  | 003145-76-4 | 87   |
| 4    |    |   | Indan, 1-methyl-                 | 132 | C10H12  | 000767-58-8 | 87   |
| 5    |    |   | 1-Phenyl-1-butene                | 132 | C10H12  | 000824-90-8 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

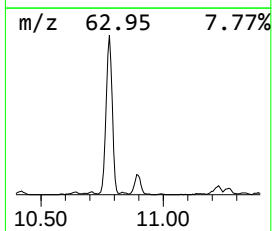
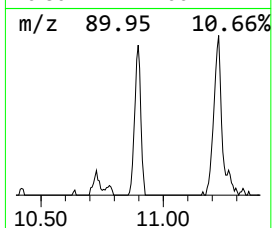
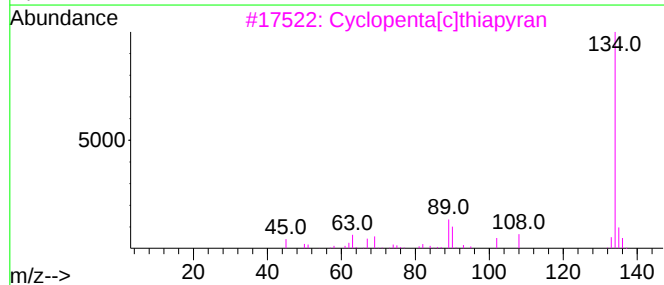
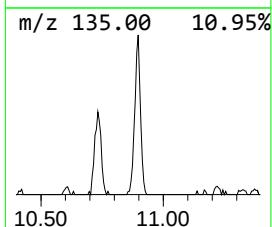
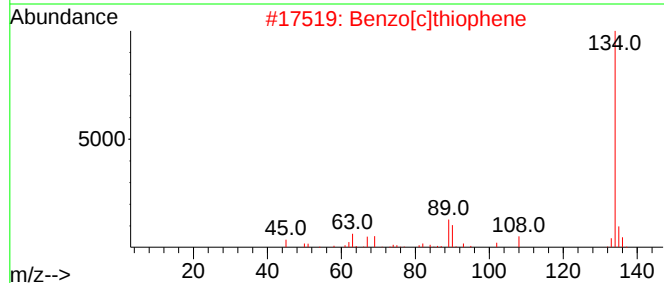
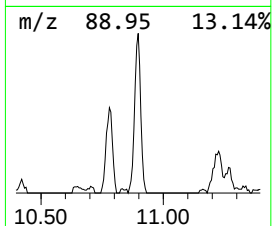
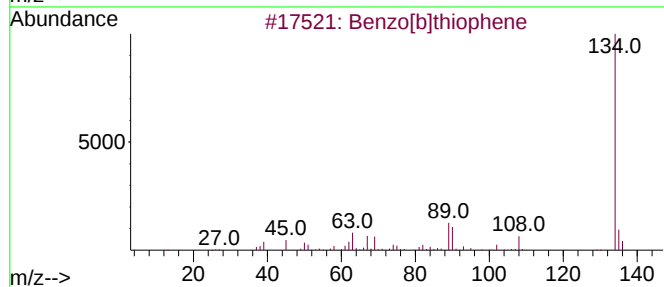
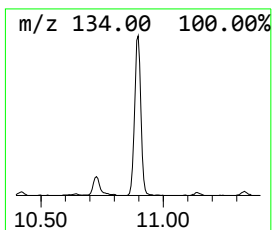
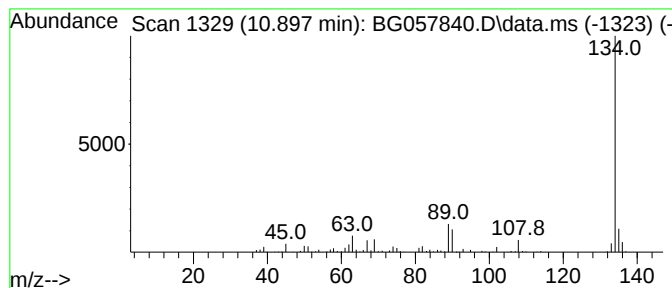
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 18 1H,1'H-2,2'-Biimidazole Concentration Rank 18

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 10.897 | 13.47 ng | 274891 | Naphthalene-d8   | 10.727 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Benzo[b]thiophene                   | 134 | C8H6S      | 000095-15-8 | 97   |
| 2    |    |   | Benzo[c]thiophene                   | 134 | C8H6S      | 000270-82-6 | 97   |
| 3    |    |   | Cyclopenta[c]thiapyran              | 134 | C8H6S      | 000270-63-3 | 94   |
| 4    |    |   | Thiazolidin-4-one, 5-benzylideno... | 287 | C13H9N3OS2 | 351062-07-2 | 50   |
| 5    |    |   | 1H,1'H-2,2'-Biimidazole             | 134 | C6H6N4     | 000492-98-8 | 42   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

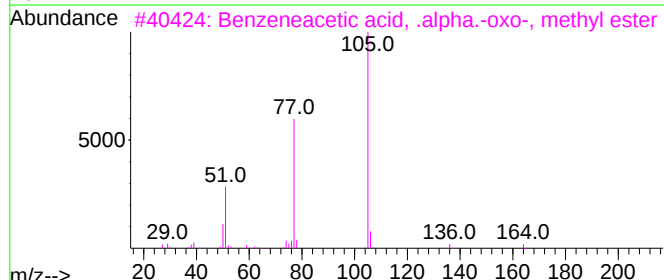
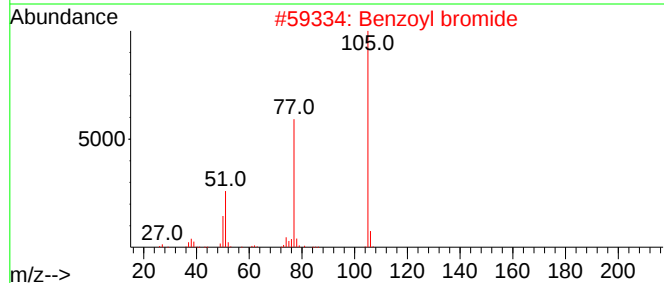
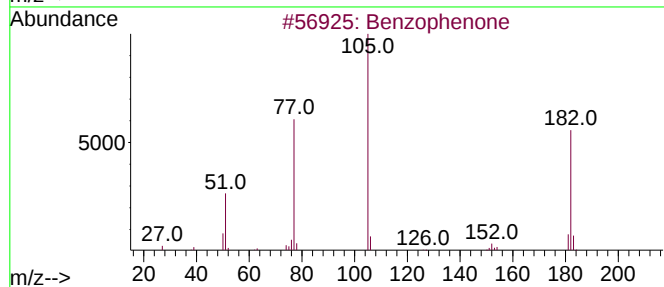
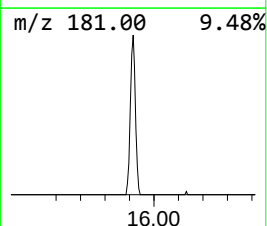
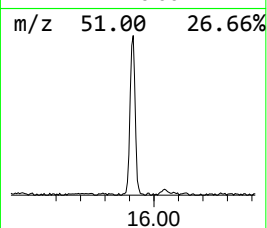
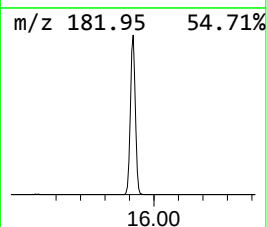
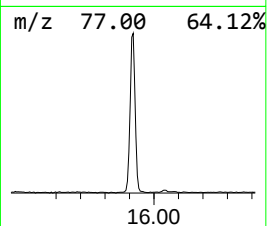
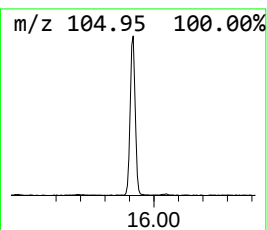
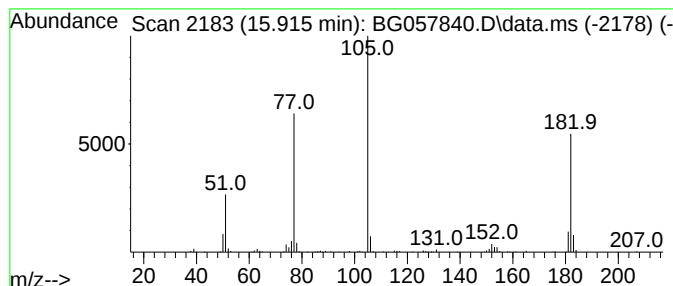
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 Benzophenone Concentration Rank 17

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 15.915 | 14.11 ng | 383565 | Acenaphthene-d10 | 14.557 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Benzophenone                        | 182 | C13H10O  | 000119-61-9 | 96   |
| 2    |    |   | Benzoyl bromide                     | 184 | C7H5BrO  | 000618-32-6 | 47   |
| 3    |    |   | Benzeneacetic acid, .alpha.-oxo-... | 164 | C9H8O3   | 015206-55-0 | 47   |
| 4    |    |   | 2-Pyrazoline-3-carbonitrile, 1-b... | 199 | C11H9N3O | 067872-68-8 | 47   |
| 5    |    |   | Benzenebutanoic acid, .gamma.-oxo-  | 178 | C10H10O3 | 002051-95-8 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062323\  
Data File : BG057840.D  
Acq On : 23 Jun 2023 13:22  
Operator : CG/JU  
Sample : 03291-03  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
MW-6A

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG061923.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 |
| Benzene, 1-ethy... | 7.013  | 221.8   | ng    | 2805990  | 1                     | 7.924  | 253083 | 20.0 |
| Benzene, 1-ethy... | 7.313  | 110.7   | ng    | 1400410  | 1                     | 7.924  | 253083 | 20.0 |
| Benzene, 1,2,3-... | 7.577  | 373.6   | ng    | 4728190  | 1                     | 7.924  | 253083 | 20.0 |
| Mesitylene         | 8.036  | 124.4   | ng    | 1574660  | 1                     | 7.924  | 253083 | 20.0 |
| Indane             | 8.300  | 128.6   | ng    | 1626940  | 1                     | 7.924  | 253083 | 20.0 |
| Indene             | 8.465  | 32.4    | ng    | 410524   | 1                     | 7.924  | 253083 | 20.0 |
| Benzene, 4-ethy... | 8.559  | 35.4    | ng    | 447744   | 1                     | 7.924  | 253083 | 20.0 |
| Benzene, 2-ethy... | 9.029  | 38.5    | ng    | 487014   | 1                     | 7.924  | 253083 | 20.0 |
| Benzene, 1,2,4,... | 9.552  | 11.9    | ng    | 242424   | 2                     | 10.727 | 408130 | 20.0 |
| Benzene, 1,2,3,... | 9.610  | 15.8    | ng    | 322645   | 2                     | 10.727 | 408130 | 20.0 |
| Phenol, 2-ethyl-   | 9.687  | 20.9    | ng    | 426435   | 2                     | 10.727 | 408130 | 20.0 |
| Benzene, 1-meth... | 9.969  | 19.9    | ng    | 405840   | 2                     | 10.727 | 408130 | 20.0 |
| Benzene, 2-ethe... | 10.121 | 36.9    | ng    | 752376   | 2                     | 10.727 | 408130 | 20.0 |
| 1H,1'H-2,2'-Bii... | 10.897 | 13.5    | ng    | 274891   | 2                     | 10.727 | 408130 | 20.0 |
| Benzophenone       | 15.915 | 14.1    | ng    | 383565   | 3                     | 14.557 | 543587 | 20.0 |