

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
 Data File : BF122590.D  
 Acq On : 8 Dec 2020 16:10  
 Operator : JU/CG  
 Sample : L4970-12  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 4S1-A-D-COMP

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\_F\METHODS\8270-BF120520.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF122590.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         | 2.216       | 18            | 22          | 31           | rVB      | 103267         | 153648        | 4.09%           | 0.505%        |
| 2         | 2.634       | 88            | 93          | 97           | rVB      | 49632          | 75429         | 2.01%           | 0.248%        |
| 3         | 2.687       | 97            | 102         | 110          | rVB2     | 38547          | 63697         | 1.70%           | 0.209%        |
| 4         | 4.152       | 347           | 351         | 358          | rVB      | 45951          | 63193         | 1.68%           | 0.208%        |
| 5         | 4.416       | 392           | 396         | 402          | rBV      | 404361         | 470698        | 12.53%          | 1.546%        |
| 6         | 4.710       | 442           | 446         | 454          | rVB4     | 50937          | 91112         | 2.43%           | 0.299%        |
| 7         | 5.046       | 498           | 503         | 509          | rVB      | 520742         | 634081        | 16.88%          | 2.083%        |
| 8         | 5.457       | 568           | 573         | 576          | rBV      | 3290122        | 3146644       | 83.78%          | 10.336%       |
| 9         | 5.851       | 636           | 640         | 645          | rBV      | 88670          | 87269         | 2.32%           | 0.287%        |
| 10        | 6.469       | 740           | 745         | 750          | rBV      | 3243361        | 3323017       | 88.48%          | 10.915%       |
| 11        | 6.651       | 770           | 776         | 778          | rBV      | 177568         | 141552        | 3.77%           | 0.465%        |
| 12        | 6.840       | 804           | 808         | 813          | rVB      | 1018878        | 893094        | 23.78%          | 2.934%        |
| 13        | 6.898       | 813           | 818         | 821          | rVV2     | 73927          | 103241        | 2.75%           | 0.339%        |
| 14        | 6.951       | 824           | 827         | 833          | rVB3     | 56997          | 57611         | 1.53%           | 0.189%        |
| 15        | 7.398       | 899           | 903         | 906          | rBV      | 2179478        | 2005114       | 53.39%          | 6.586%        |
| 16        | 7.545       | 925           | 928         | 933          | rVB3     | 67266          | 89122         | 2.37%           | 0.293%        |
| 17        | 7.681       | 944           | 951         | 955          | rVB3     | 53360          | 81416         | 2.17%           | 0.267%        |
| 18        | 8.122       | 1022          | 1026        | 1029         | rBV      | 1484851        | 1209206       | 32.20%          | 3.972%        |
| 19        | 8.163       | 1029          | 1033        | 1038         | rVB2     | 52633          | 71988         | 1.92%           | 0.236%        |
| 20        | 8.698       | 1121          | 1124        | 1128         | rBV6     | 41407          | 57611         | 1.53%           | 0.189%        |
| 21        | 9.198       | 1205          | 1209        | 1212         | rBV      | 4474427        | 3755655       | 100.00%         | 12.336%       |
| 22        | 9.592       | 1272          | 1276        | 1282         | rBV      | 307609         | 321375        | 8.56%           | 1.056%        |
| 23        | 9.875       | 1321          | 1324        | 1328         | rBV      | 1662741        | 1419820       | 37.80%          | 4.664%        |
| 24        | 10.528      | 1433          | 1435        | 1438         | rVB2     | 165881         | 159801        | 4.25%           | 0.525%        |
| 25        | 10.663      | 1454          | 1458        | 1466         | rBV      | 2295645        | 2467493       | 65.70%          | 8.105%        |
| 26        | 11.257      | 1557          | 1559        | 1563         | rBV      | 211615         | 182168        | 4.85%           | 0.598%        |
| 27        | 11.363      | 1574          | 1577        | 1580         | rBV      | 1509384        | 1179716       | 31.41%          | 3.875%        |
| 28        | 11.492      | 1597          | 1599        | 1601         | rBV2     | 299282         | 224270        | 5.97%           | 0.737%        |
| 29        | 11.522      | 1602          | 1604        | 1606         | rBV      | 273362         | 266479        | 7.10%           | 0.875%        |
| 30        | 11.592      | 1614          | 1616        | 1621         | rVB3     | 280384         | 309434        | 8.24%           | 1.016%        |
| 31        | 11.757      | 1642          | 1644        | 1647         | rBV3     | 283277         | 329228        | 8.77%           | 1.081%        |
| 32        | 11.904      | 1667          | 1669        | 1673         | rVB2     | 596861         | 558283        | 14.87%          | 1.834%        |
| 33        | 12.698      | 1801          | 1804        | 1807         | rVB      | 1840961        | 1657955       | 44.15%          | 5.446%        |
| 34        | 12.957      | 1844          | 1848        | 1851         | rVB      | 3271648        | 3032052       | 80.73%          | 9.959%        |
| 35        | 14.010      | 2024          | 2027        | 2029         | rVB      | 985608         | 913528        | 24.32%          | 3.001%        |
| 36        | 15.468      | 2272          | 2275        | 2279         | rBV      | 734564         | 848025        | 22.58%          | 2.786%        |

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Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

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\_F\METHODS\8270-BF120520.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

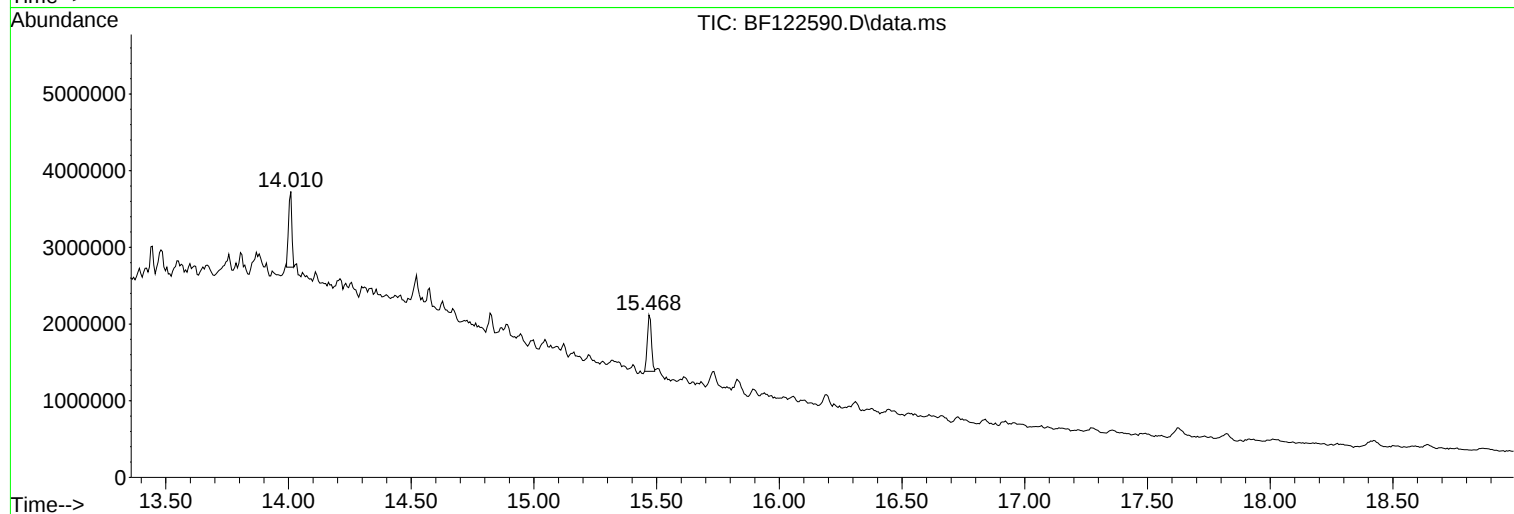
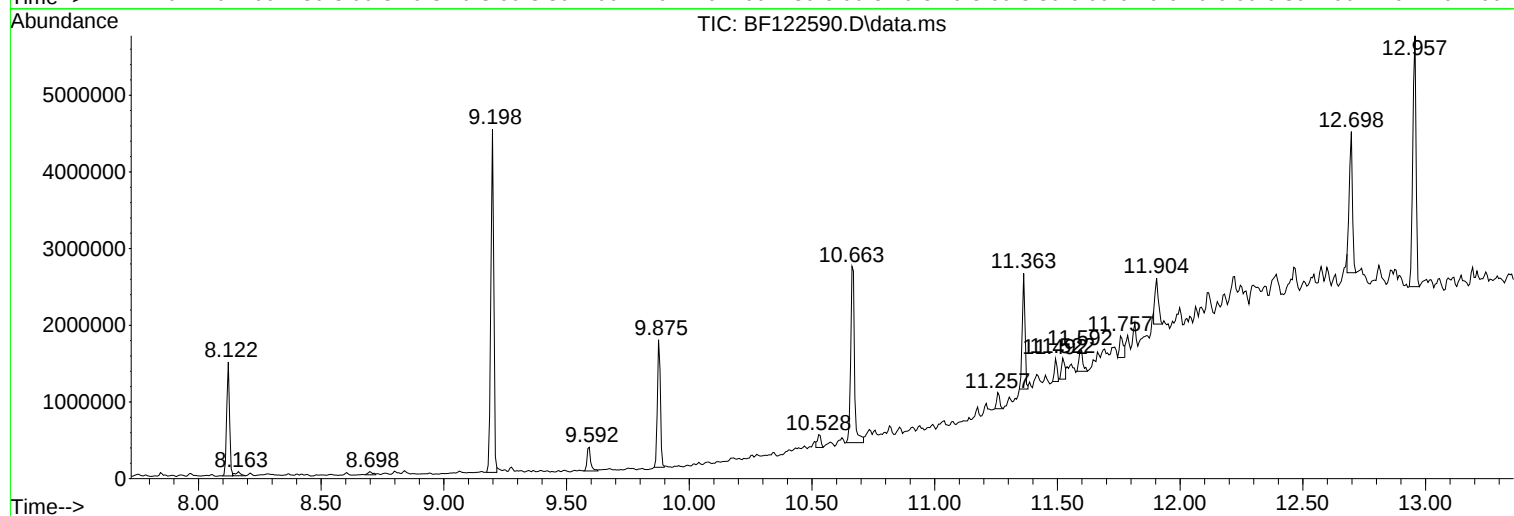
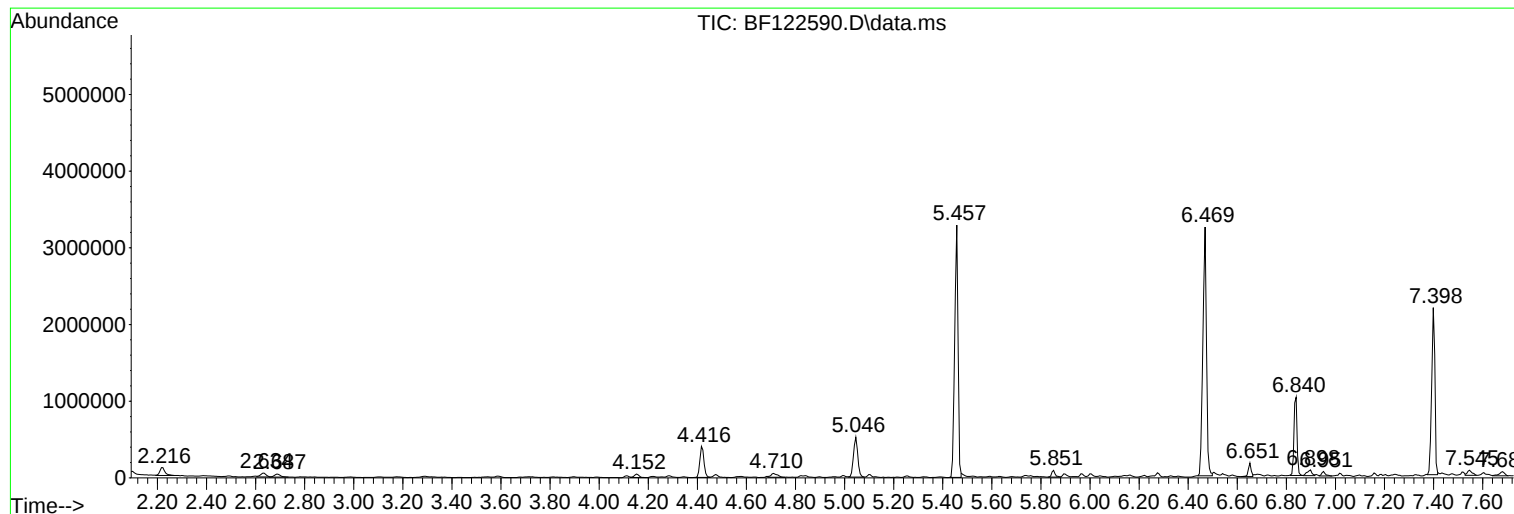
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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ClientSampleId :  
4S1-A-D-COMP

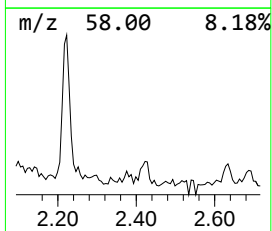
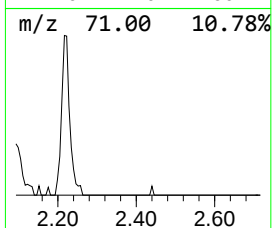
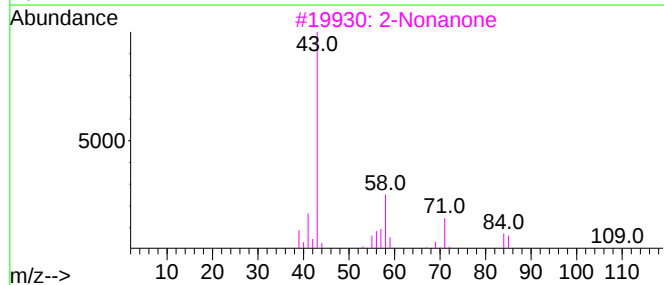
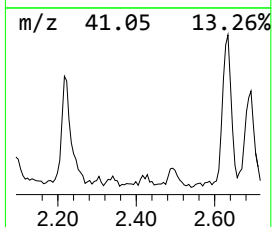
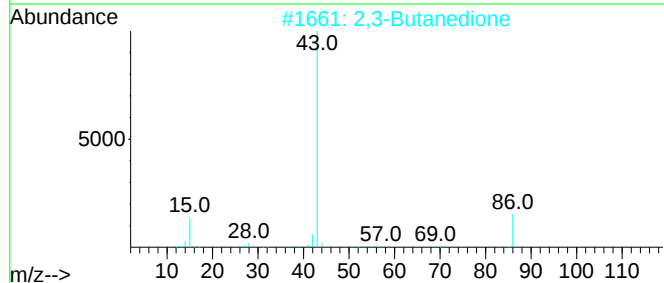
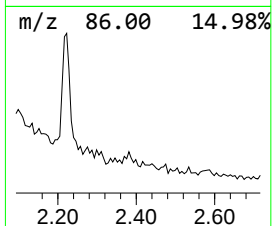
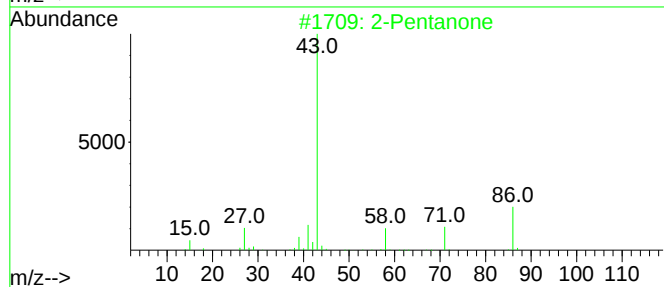
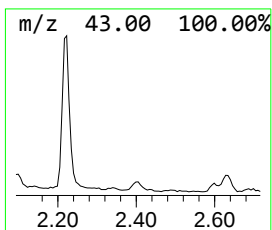
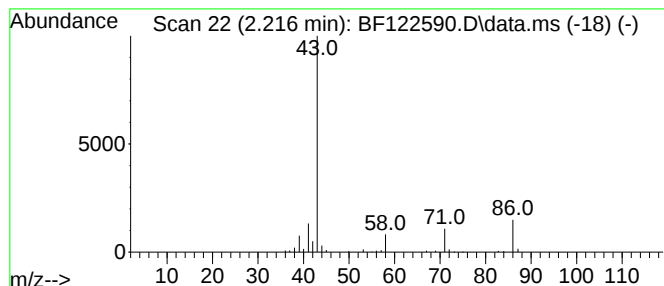
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 2-Pentanone Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 2.216 | 3.44 ng | 153648 | 1,4-Dichlorobenzene-d4 | 6.840 |

| Hit# | of 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------|-----|---------|-------------|------|
| 1    |      | 2-Pentanone           | 86  | C5H10O  | 000107-87-9 | 80   |
| 2    |      | 2,3-Butanedione       | 86  | C4H6O2  | 000431-03-8 | 47   |
| 3    |      | 2-Nonanone            | 142 | C9H18O  | 000821-55-6 | 33   |
| 4    |      | 2-Butanone, 3-methyl- | 86  | C5H10O  | 000563-80-4 | 23   |
| 5    |      | 2,3-Hexanedione       | 114 | C6H10O2 | 003848-24-6 | 17   |



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BNA\_F  
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4S1-A-D-COMP

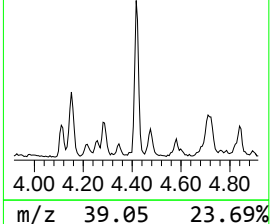
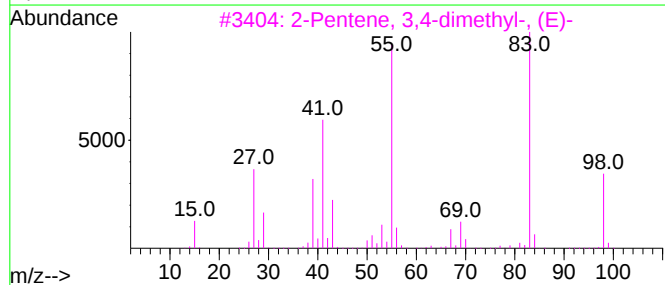
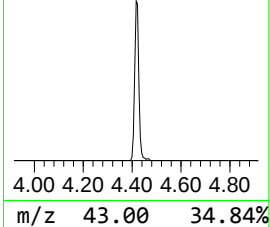
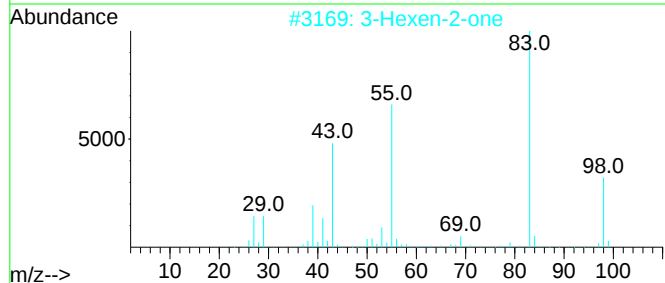
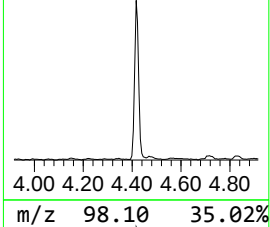
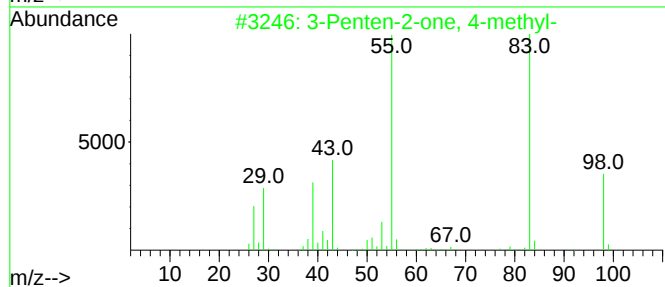
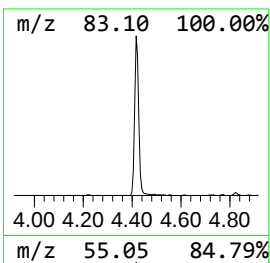
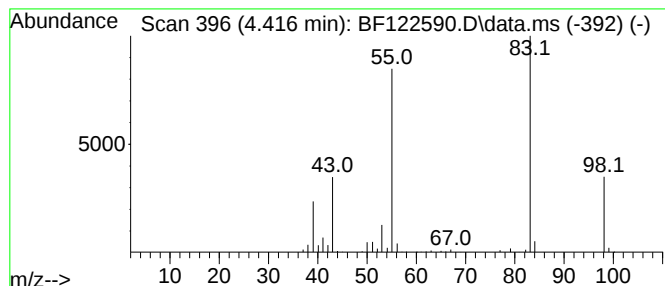
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 4.416 | 10.54 ng | 470698 | 1,4-Dichlorobenzene-d4 | 6.840 |

| Hit# | of 5 | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|------|--------------------------------|----|---------|-------------|------|
| 1    |      | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 95   |
| 2    |      | 3-Hexen-2-one                  | 98 | C6H10O  | 000763-93-9 | 91   |
| 3    |      | 2-Pentene, 3,4-dimethyl-, (E)- | 98 | C7H14   | 004914-92-5 | 90   |
| 4    |      | 2-Pentene, 3,4-dimethyl-, (Z)- | 98 | C7H14   | 004914-91-4 | 86   |
| 5    |      | 2-Pentene, 2,4-dimethyl-       | 98 | C7H14   | 000625-65-0 | 80   |



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Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

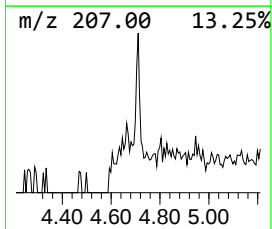
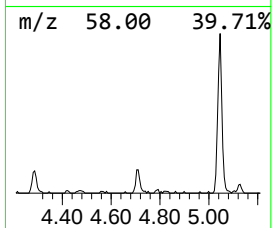
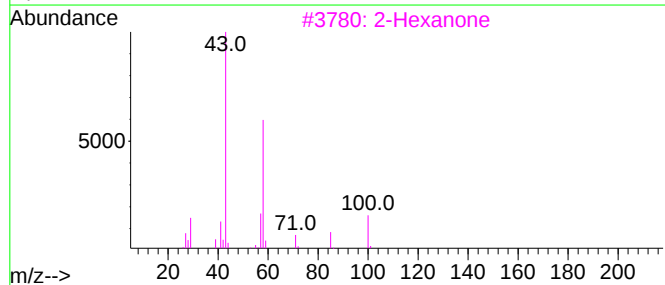
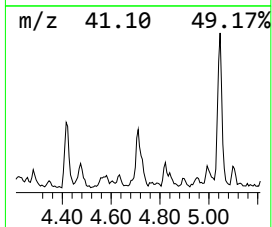
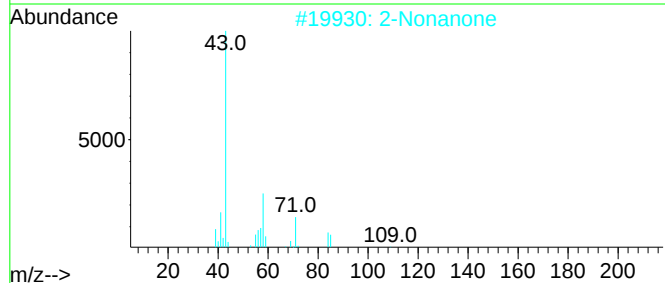
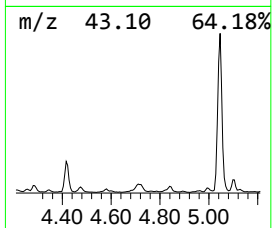
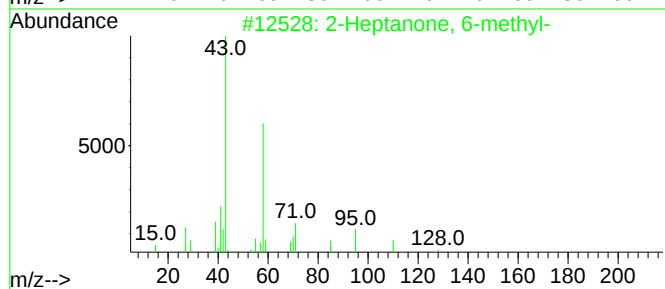
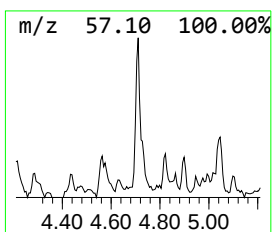
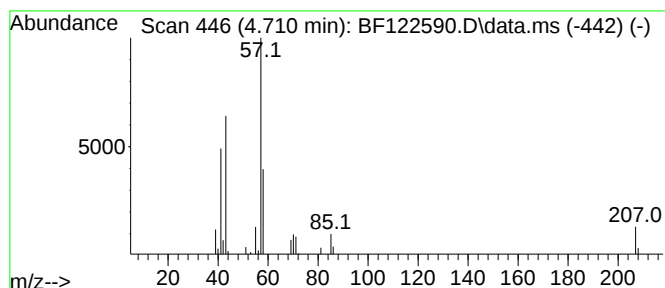
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown4.71 Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD       | R.T.  |
|-------|---------|-------|------------------------|-------|
| 4.710 | 2.04 ng | 91112 | 1,4-Dichlorobenzene-d4 | 6.840 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Heptanone, 6-methyl- | 128 | C8H16O  | 000928-68-7 | 28   |
| 2    |    |   | 2-Nonanone             | 142 | C9H18O  | 000821-55-6 | 25   |
| 3    |    |   | 2-Hexanone             | 100 | C6H12O  | 000591-78-6 | 25   |
| 4    |    |   | 2-Hexanone, 4-methyl-  | 114 | C7H14O  | 000105-42-0 | 23   |
| 5    |    |   | 2-Hexanone, 5-methyl-  | 114 | C7H14O  | 000110-12-3 | 23   |



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Instrument :  
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4S1-A-D-COMP

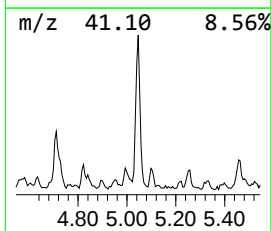
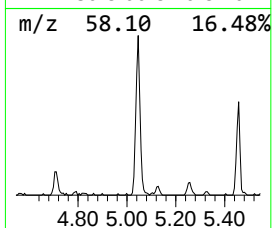
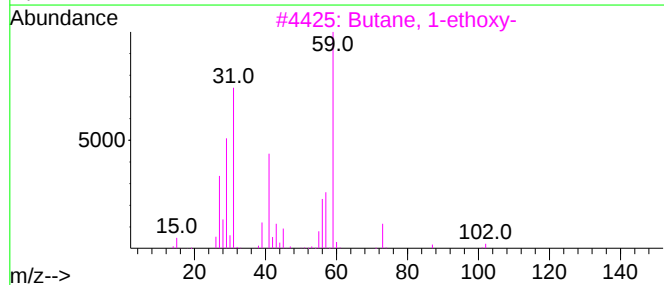
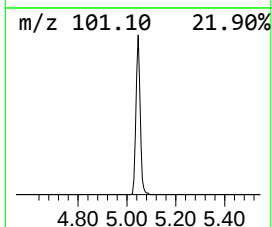
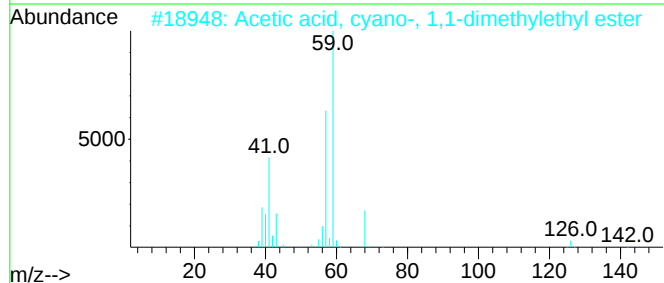
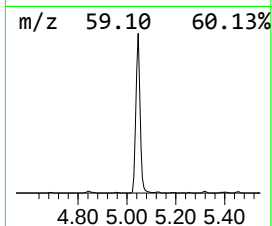
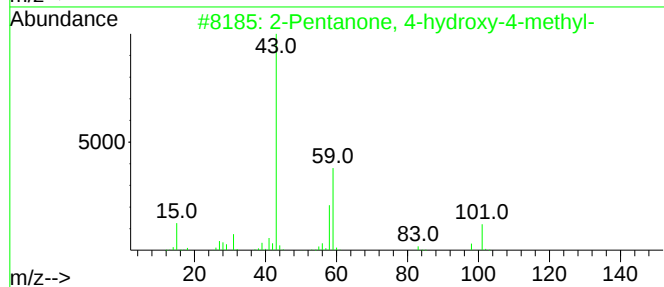
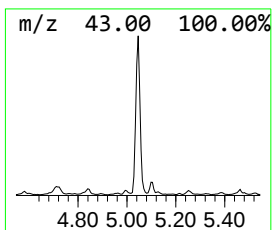
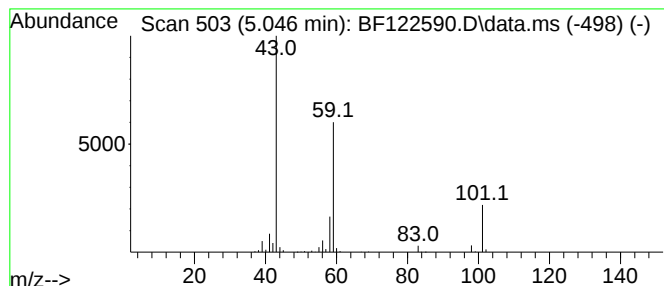
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.      | EstConc                             | Area   | Relative to ISTD       | R.T.        |      |
|-----------|-------------------------------------|--------|------------------------|-------------|------|
| 5.046     | 14.20 ng                            | 634081 | 1,4-Dichlorobenzene-d4 | 6.840       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm                | CAS#        | Qual |
| 1         | 2-Pentanone, 4-hydroxy-4-methyl-    | 116    | C6H12O2                | 000123-42-2 | 50   |
| 2         | Acetic acid, cyano-, 1,1-dimethy... | 141    | C7H11NO2               | 001116-98-9 | 32   |
| 3         | Butane, 1-ethoxy-                   | 102    | C6H14O                 | 000628-81-9 | 9    |
| 4         | Morpholine, 4-methyl-               | 101    | C5H11NO                | 000109-02-4 | 9    |
| 5         | (+)-4-Amino-4,5-dihydro-2(3H)-f...  | 101    | C4H7NO2                | 016504-58-8 | 9    |



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Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

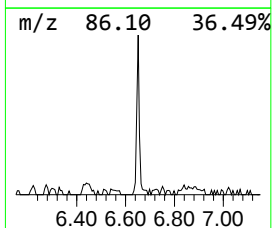
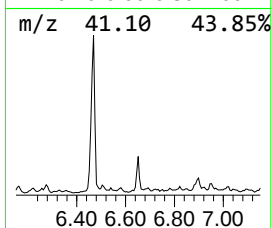
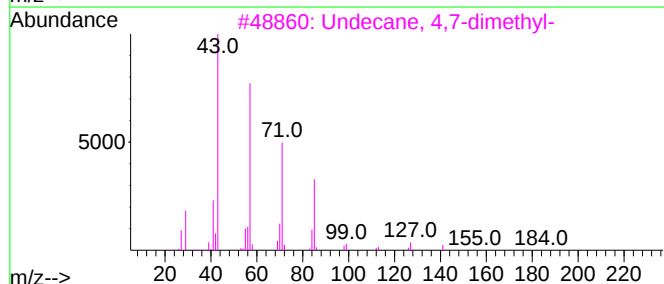
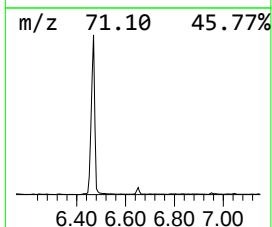
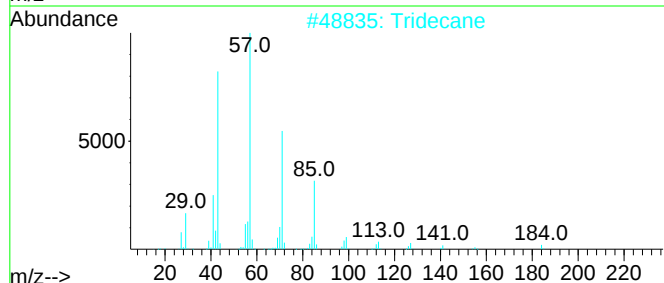
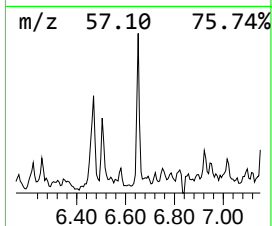
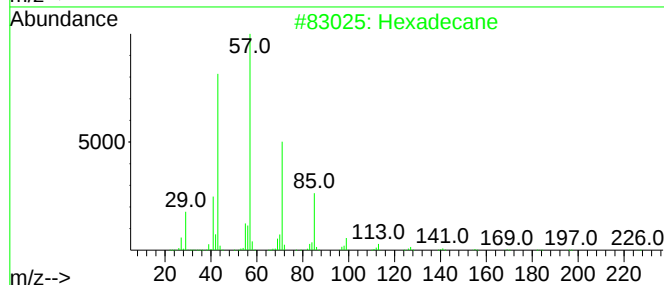
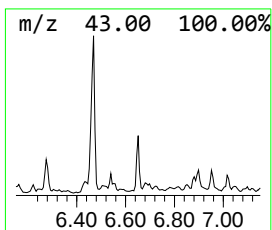
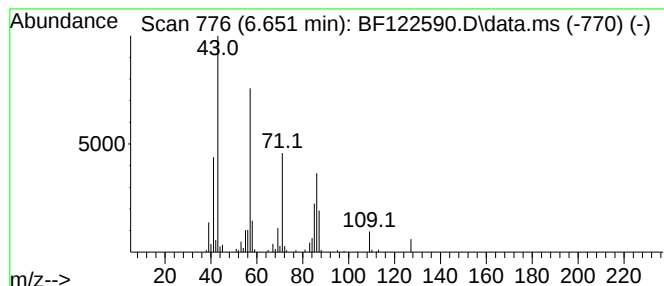
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 unknown6.65 Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.651 | 3.17 ng | 141552 | 1,4-Dichlorobenzene-d4 | 6.840 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 47   |
| 2    |    |   | Tridecane                           | 184 | C13H28   | 000629-50-5  | 43   |
| 3    |    |   | Undecane, 4,7-dimethyl-             | 184 | C13H28   | 017301-32-5  | 38   |
| 4    |    |   | Decane, 2,4-dimethyl-               | 170 | C12H26   | 002801-84-5  | 38   |
| 5    |    |   | Oxalic acid, 2-ethylhexyl hexyl ... | 286 | C16H30O4 | 1000309-38-9 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

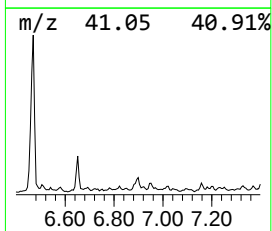
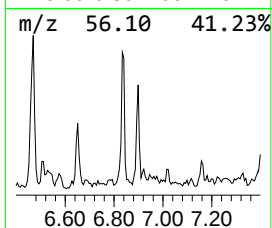
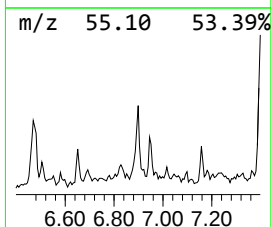
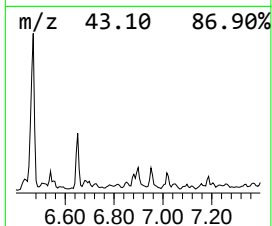
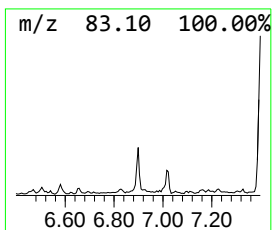
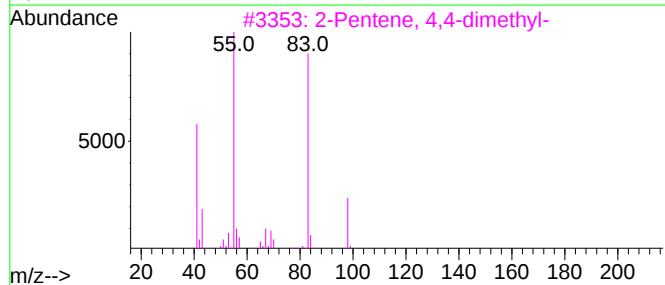
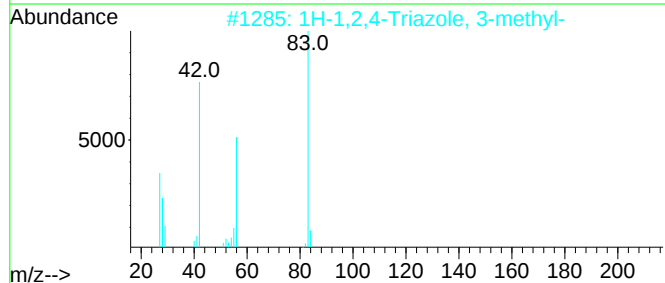
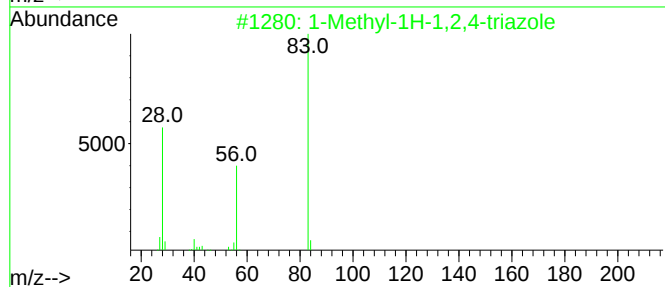
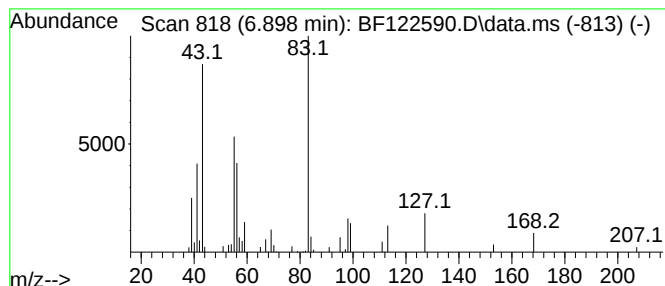
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Peak Number 6 1-Methyl-1H-1,2,4-triazole Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.898 | 2.31 ng | 103241 | 1,4-Dichlorobenzene-d4 | 6.840 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 1-Methyl-1H-1,2,4-triazole          | 83  | C3H5N3  | 006086-21-1 | 52   |
| 2    |      | 1H-1,2,4-Triazole, 3-methyl-        | 83  | C3H5N3  | 007170-01-6 | 50   |
| 3    |      | 2-Pentene, 4,4-dimethyl-            | 98  | C7H14   | 026232-98-4 | 43   |
| 4    |      | 2-Pentene, 4,4-dimethyl-, (E)-      | 98  | C7H14   | 000690-08-4 | 43   |
| 5    |      | 3-Hexene, 2,5-dimethyl-3,4-bis(1... | 196 | C14H28  | 007090-88-2 | 37   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

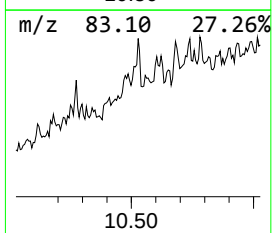
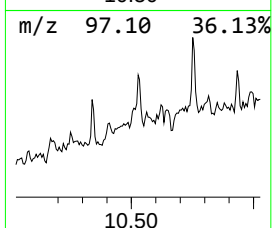
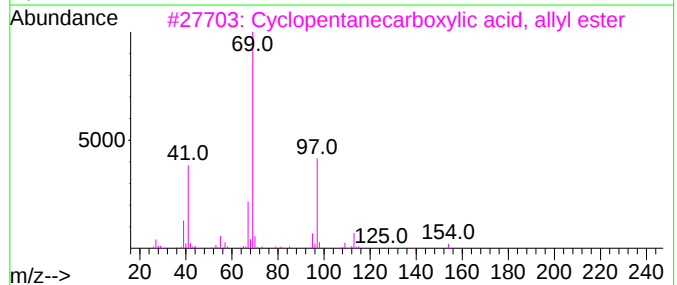
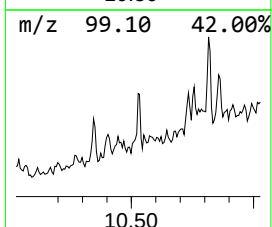
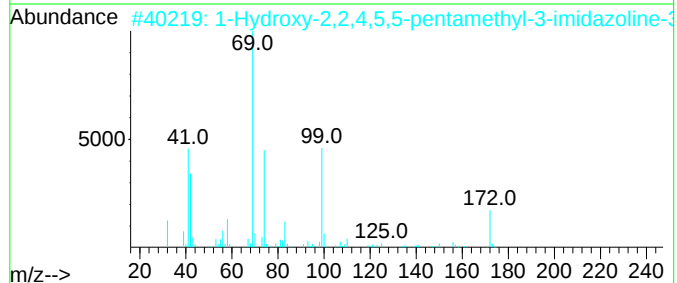
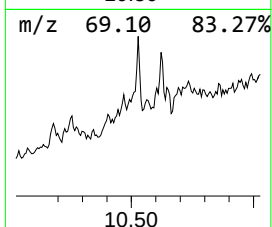
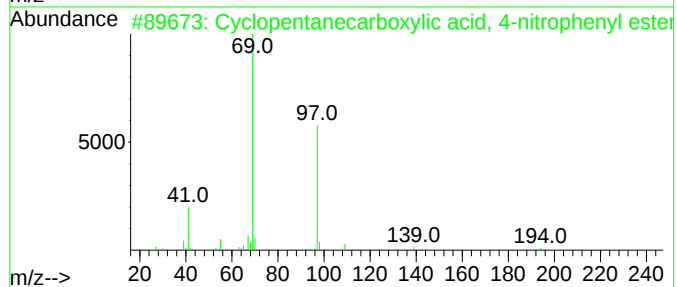
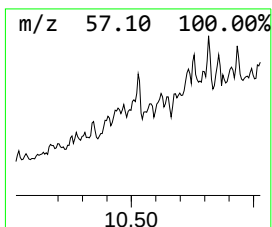
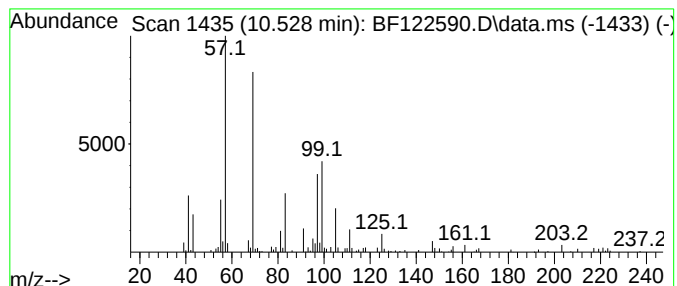
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 unknown10.52 Concentration Rank 13

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.528    | 2.25 ng                             | 159801 | Acenaphthene-d10 | 9.875        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Cyclopentanecarboxylic acid, 4-n... | 235    | C12H13NO4        | 1000307-58-0 | 35   |
| 2         | 1-Hydroxy-2,2,4,5,5-pentamethyl-... | 172    | C8H16N2O2        | 018796-02-6  | 28   |
| 3         | Cyclopentanecarboxylic acid, all... | 154    | C9H14O2          | 178905-33-4  | 25   |
| 4         | 1-Nitrododecane                     | 215    | C12H25NO2        | 016891-99-9  | 20   |
| 5         | 2-Propionyl-1-pyrroline             | 125    | C7H11NO          | 1000298-55-4 | 17   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

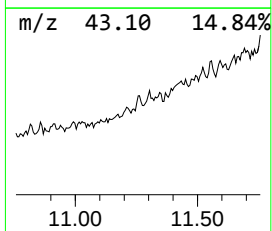
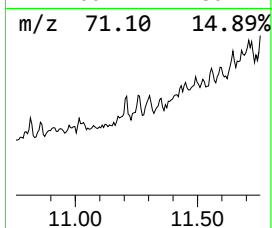
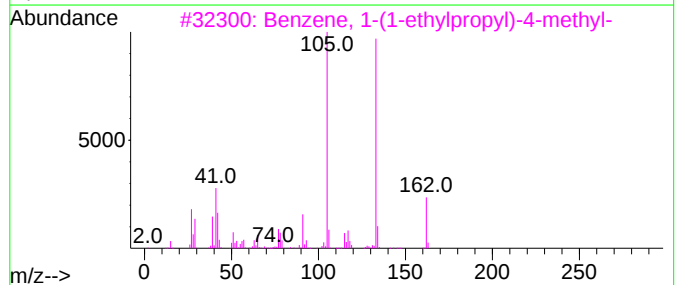
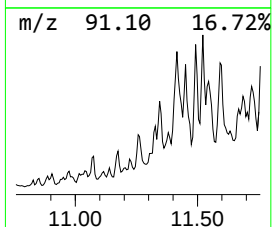
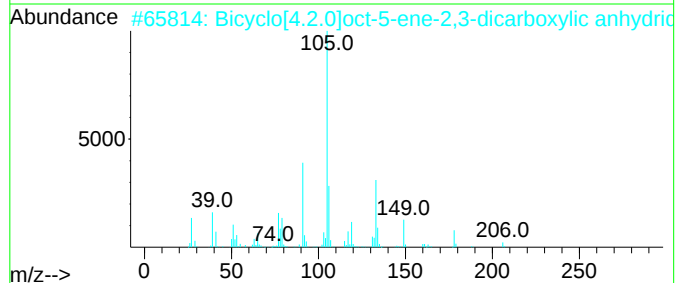
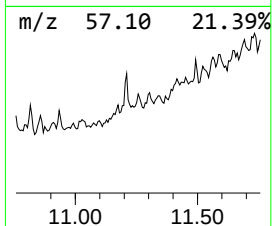
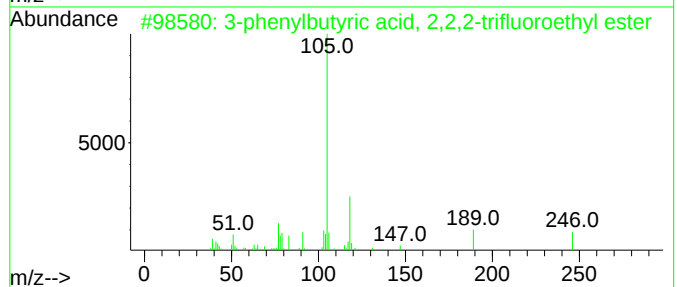
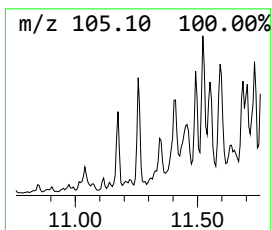
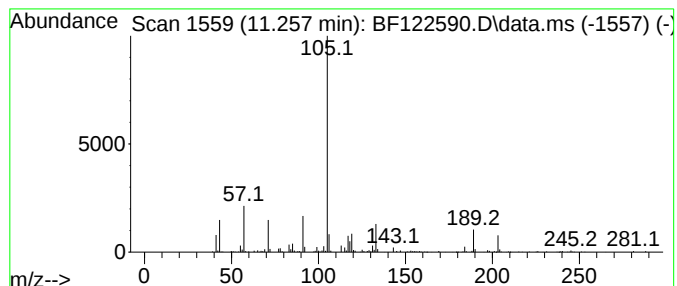
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown11.25 Concentration Rank 11

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.257 | 3.09 ng | 182168 | Phenanthrene-d10 | 11.363 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 3-phenylbutyric acid, 2,2,2-trif... | 246 | C12H13F3O2 | 1000376-55-4 | 42   |
| 2    |    |   | Bicyclo[4.2.0]oct-5-ene-2,3-dica... | 206 | C12H14O3   | 022467-69-2  | 33   |
| 3    |    |   | Benzene, 1-(1-ethylpropyl)-4-met... | 162 | C12H18     | 022975-58-2  | 28   |
| 4    |    |   | 3-Cyanophenyl-.beta.-phenylpropi... | 251 | C16H13NO2  | 040123-39-5  | 27   |
| 5    |    |   | 4,5-Dihydro-4,4-dimethyl-2-pheny... | 203 | C12H13NO2  | 029637-55-6  | 16   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

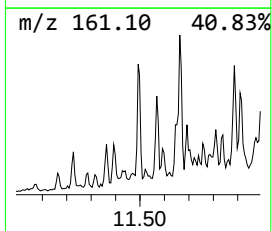
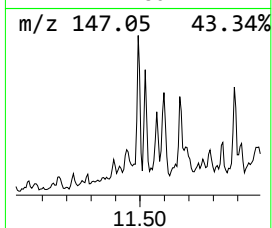
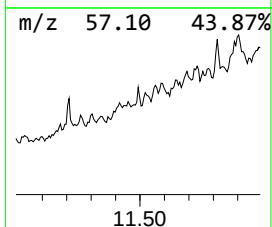
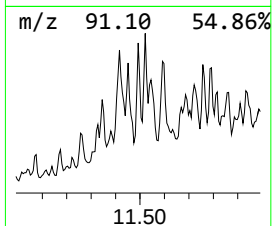
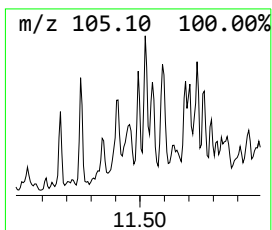
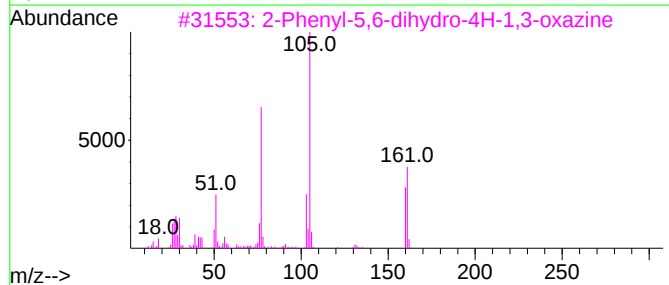
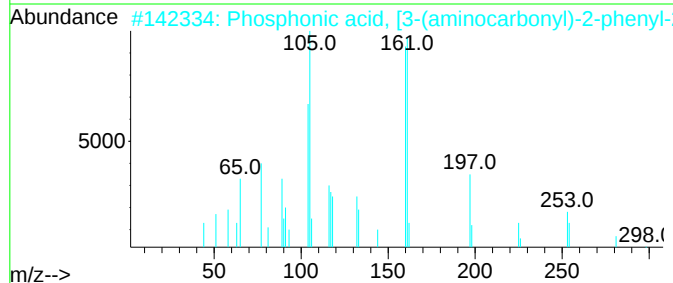
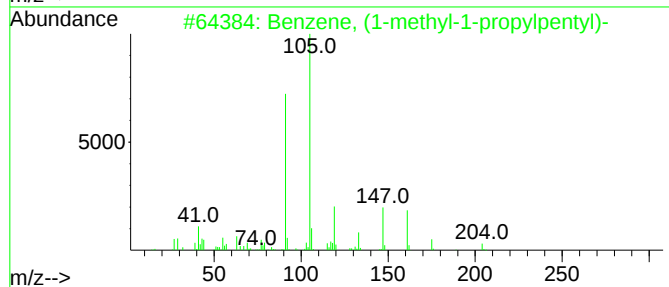
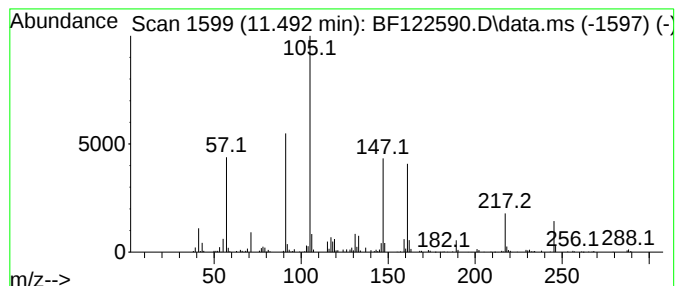
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 unknown11.49 Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.492 | 3.80 ng | 224270 | Phenanthrene-d10 | 11.363 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | Benzene, (1-methyl-1-propylpentyl)- | 204 | C15H24      | 054932-91-1 | 38   |
| 2    |    |   | Phosphonic acid, [3-(aminocarbon... | 298 | C13H19N2O4P | 033866-52-3 | 27   |
| 3    |    |   | 2-Phenyl-5,6-dihydro-4H-1,3-oxazine | 161 | C10H11NO    | 019029-45-9 | 17   |
| 4    |    |   | Benzene, (1,1-diethylpropyl)-       | 176 | C13H20      | 004170-84-7 | 12   |
| 5    |    |   | .alpha.-Cubebene                    | 204 | C15H24      | 017699-14-8 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

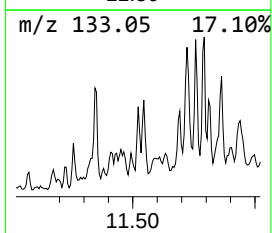
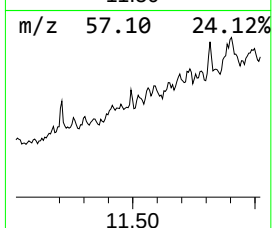
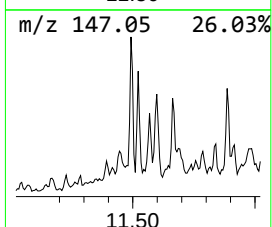
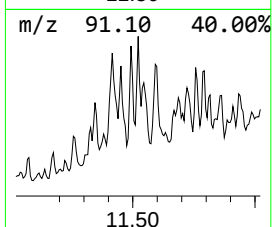
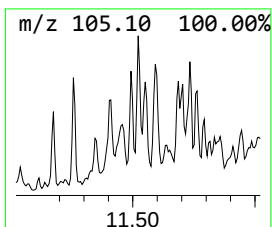
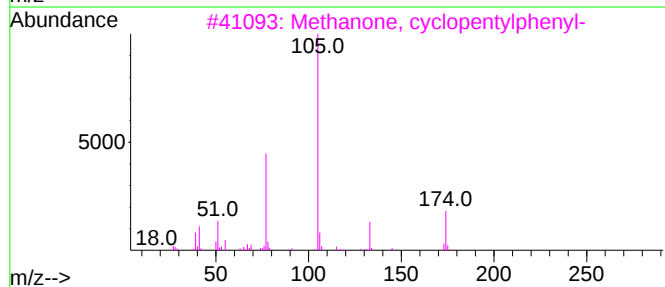
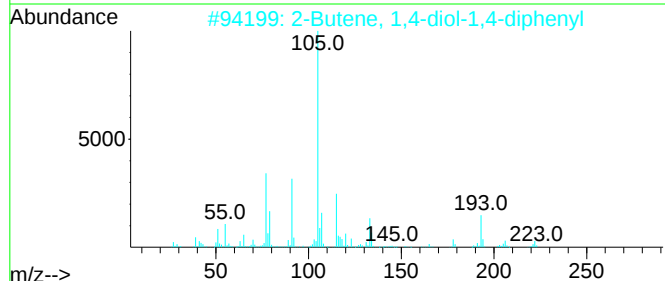
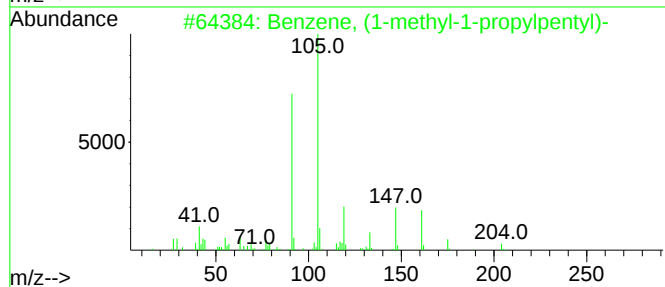
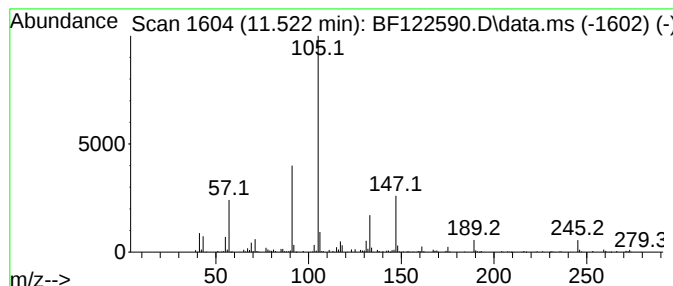
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 unknown11.52 Concentration Rank 7

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.522 | 4.52 ng | 266479 | Phenanthrene-d10 | 11.363 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|------|-------------------------------------|-----|------------|--------------|------|
| 1    |      | Benzene, (1-methyl-1-propylpentyl)- | 204 | C15H24     | 054932-91-1  | 38   |
| 2    |      | 2-Butene, 1,4-diol-1,4-diphenyl     | 240 | C16H16O2   | 1000197-44-1 | 32   |
| 3    |      | Methanone, cyclopentylphenyl-       | 174 | C12H14O    | 005422-88-8  | 25   |
| 4    |      | Malonic acid, 8-chlorooctyl prop... | 292 | C14H25ClO4 | 1000349-00-3 | 25   |
| 5    |      | Benzoic acid, 1-(2-thienyl)ethyl... | 245 | C13H11NO2S | 309282-70-0  | 25   |



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Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
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4S1-A-D-COMP

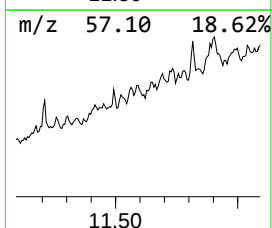
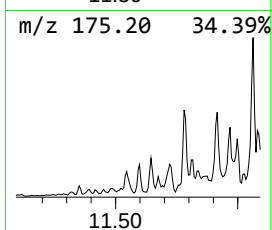
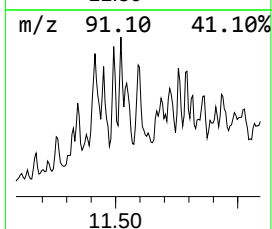
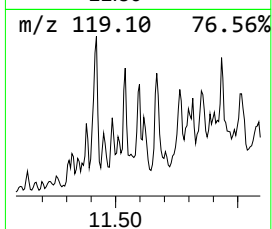
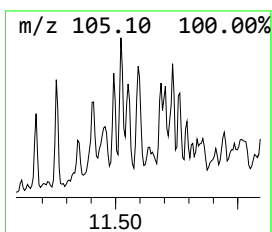
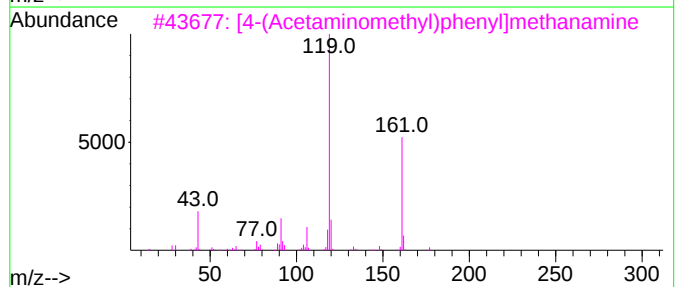
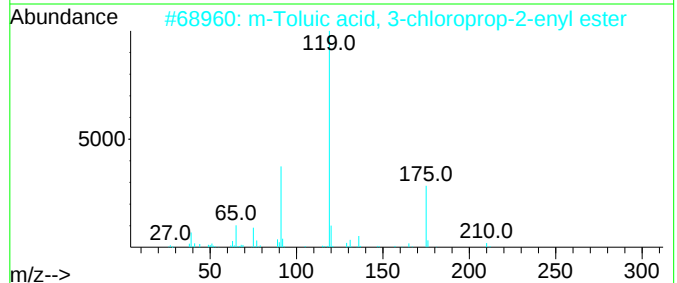
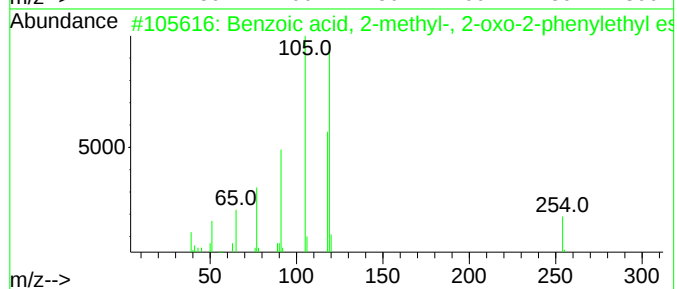
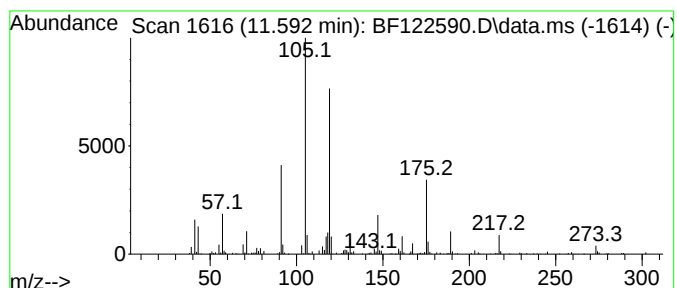
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 unknown11.59 Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 11.592    | 5.25 ng                             | 309434 | Phenanthrene-d10 | 11.363       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Benzoic acid, 2-methyl-, 2-oxo-2... | 254    | C16H14O3         | 055153-21-4  | 47   |
| 2         | m-Toluic acid, 3-chloroprop-2-en... | 210    | C11H11ClO2       | 1000292-50-0 | 47   |
| 3         | [4-(Acetaminomethyl)phenyl]metha... | 178    | C10H14N2O        | 1000373-14-1 | 46   |
| 4         | Benzonitrile, 2-ethoxy-             | 147    | C9H9NO           | 006609-57-0  | 43   |
| 5         | o-Toluic acid, 2-ethylcyclohexyl... | 246    | C16H22O2         | 1000293-34-8 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

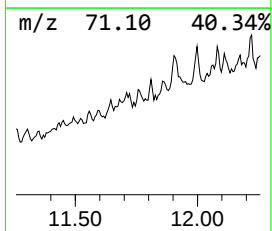
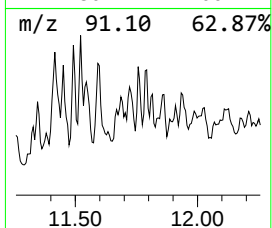
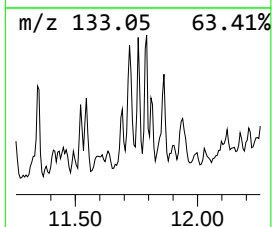
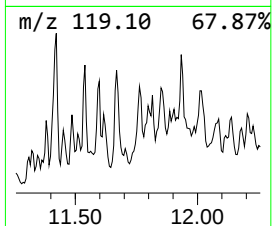
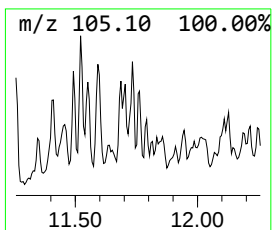
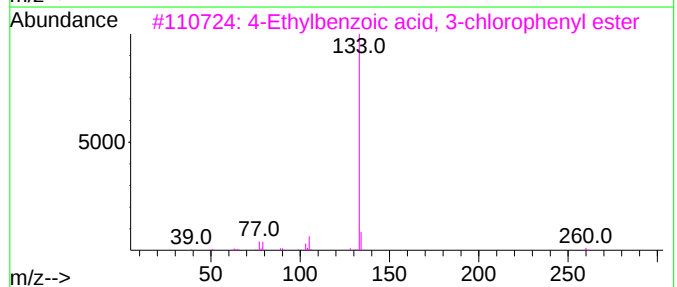
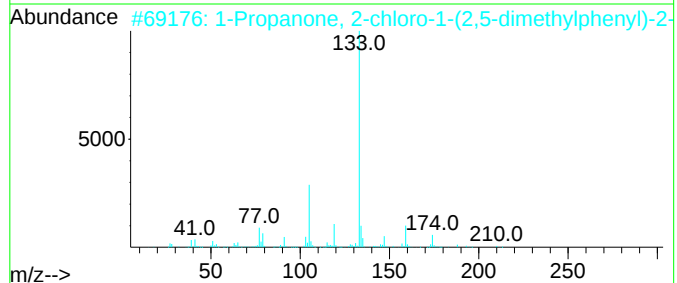
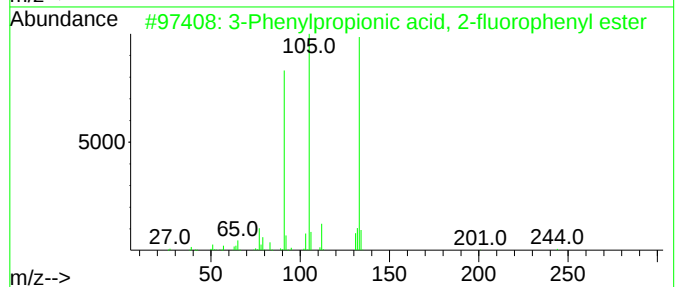
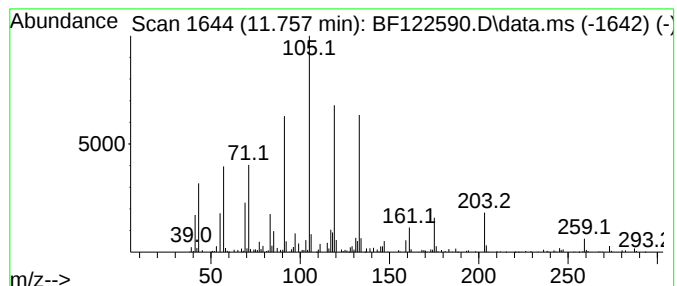
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 unknown11.75 Concentration Rank 5

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.757 | 5.58 ng | 329228 | Phenanthrene-d10 | 11.363 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|------|-------------------------------------|-----|------------|--------------|------|
| 1    |      | 3-Phenylpropionic acid, 2-fluoro... | 244 | C15H13F02  | 1000325-75-9 | 14   |
| 2    |      | 1-Propanone, 2-chloro-1-(2,5-dim... | 210 | C12H15ClO  | 054965-52-5  | 11   |
| 3    |      | 4-Ethylbenzoic acid, 3-chlorophe... | 260 | C15H13ClO2 | 1000325-73-4 | 11   |
| 4    |      | 1,2-Benzenediol, o-(4-ethylbenzo... | 242 | C15H14O3   | 1000325-96-6 | 11   |
| 5    |      | 1-Propanone, 2-chloro-1-(2,4-dim... | 210 | C12H15ClO  | 054965-53-6  | 11   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

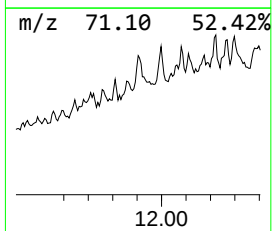
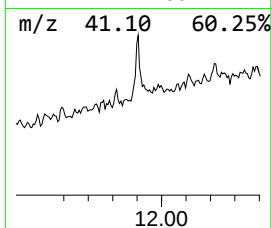
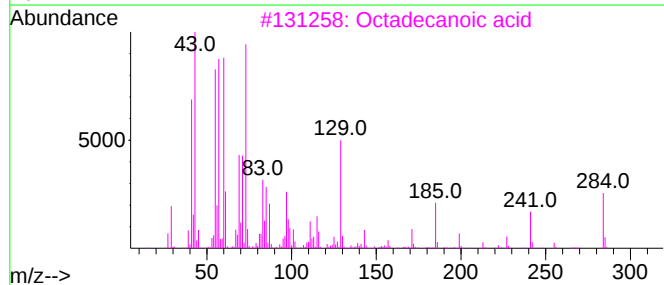
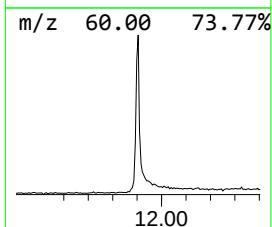
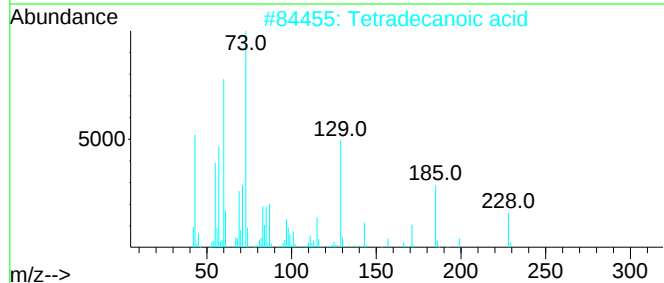
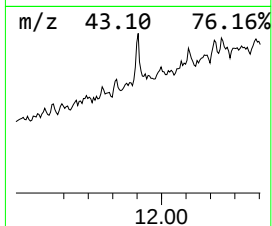
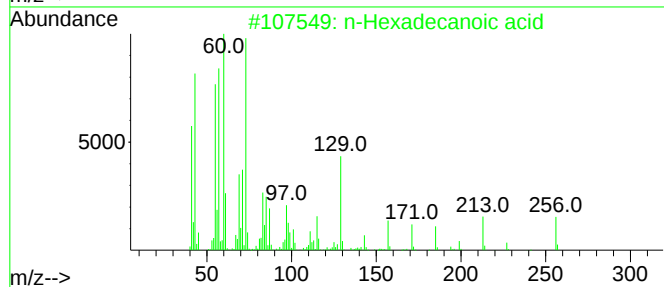
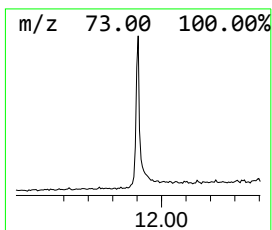
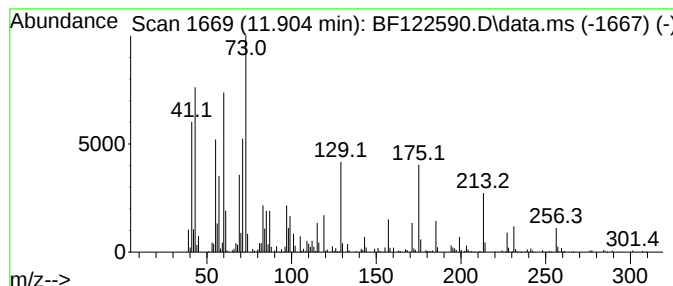
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 n-Hexadecanoic acid Concentration Rank 4

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 11.904    | 9.46 ng             | 558283 | Phenanthrene-d10 | 11.363      |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 93   |
| 2         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 86   |
| 3         | Octadecanoic acid   | 284    | C18H36O2         | 000057-11-4 | 70   |
| 4         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 53   |
| 5         | Tridecanoic acid    | 214    | C13H26O2         | 000638-53-9 | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

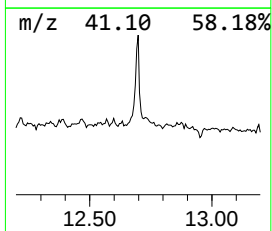
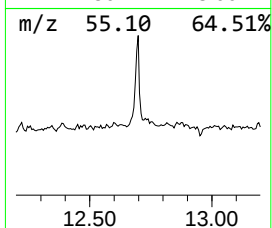
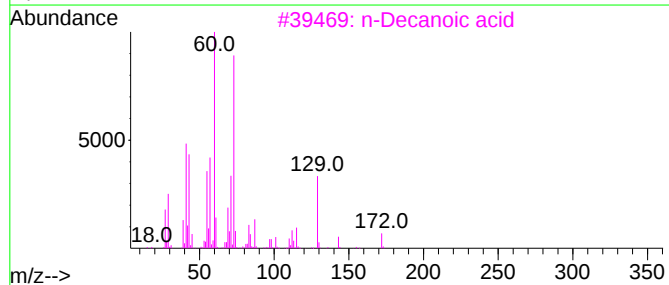
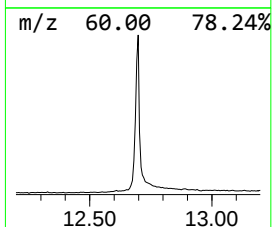
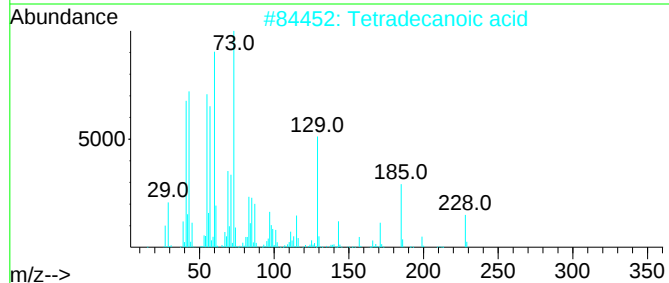
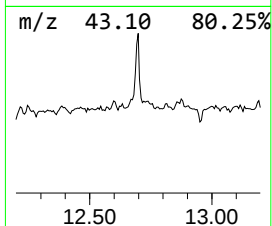
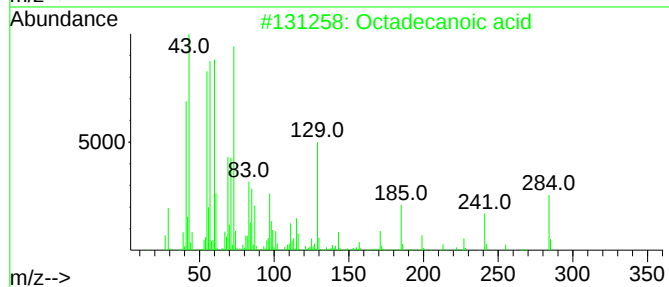
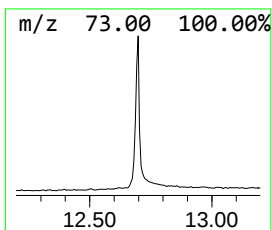
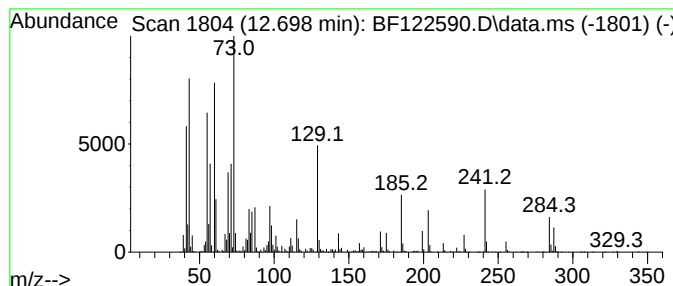
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Octadecanoic acid Concentration Rank 1

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 12.698 | 36.30 ng | 1657960 | Chrysene-d12     | 14.010 |

| Hit# | of 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|------|---------------------|-----|----------|-------------|------|
| 1    |      | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 83   |
| 2    |      | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 70   |
| 3    |      | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 60   |
| 4    |      | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 53   |
| 5    |      | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF120820\  
Data File : BF122590.D  
Acq On : 8 Dec 2020 16:10  
Operator : JU/CG  
Sample : L4970-12  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
4S1-A-D-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF120520.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| 2-Pentanone        | 2.216  | 3.4     | ng    | 153648   | 1                     | 6.840  | 893094  | 20.0 |
| 3-Penten-2-one,... | 4.416  | 10.5    | ng    | 470698   | 1                     | 6.840  | 893094  | 20.0 |
| unknown4.71        | 4.710  | 2.0     | ng    | 91112    | 1                     | 6.840  | 893094  | 20.0 |
| 2-Pentanone, 4-... | 5.046  | 14.2    | ng    | 634081   | 1                     | 6.840  | 893094  | 20.0 |
| unknown6.65        | 6.651  | 3.2     | ng    | 141552   | 1                     | 6.840  | 893094  | 20.0 |
| 1-Methyl-1H-1,2... | 6.898  | 2.3     | ng    | 103241   | 1                     | 6.840  | 893094  | 20.0 |
| unknown10.52       | 10.528 | 2.3     | ng    | 159801   | 3                     | 9.875  | 1419820 | 20.0 |
| unknown11.25       | 11.257 | 3.1     | ng    | 182168   | 4                     | 11.363 | 1179720 | 20.0 |
| unknown11.49       | 11.492 | 3.8     | ng    | 224270   | 4                     | 11.363 | 1179720 | 20.0 |
| unknown11.52       | 11.522 | 4.5     | ng    | 266479   | 4                     | 11.363 | 1179720 | 20.0 |
| unknown11.59       | 11.592 | 5.3     | ng    | 309434   | 4                     | 11.363 | 1179720 | 20.0 |
| unknown11.75       | 11.757 | 5.6     | ng    | 329228   | 4                     | 11.363 | 1179720 | 20.0 |
| n-Hexadecanoic ... | 11.904 | 9.5     | ng    | 558283   | 4                     | 11.363 | 1179720 | 20.0 |
| Octadecanoic acid  | 12.698 | 36.3    | ng    | 1657960  | 5                     | 14.010 | 913528  | 20.0 |