

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
 Data File : BF142543.D  
 Acq On : 23 May 2025 18:58  
 Operator : RC/JU  
 Sample : Q2102-03  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 LAW-25-0078

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-BF052025.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142543.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         | 4.416       | 372           | 378         | 386          | rBV      | 103602         | 158720        | 3.10%           | 0.498%        |
| 2         | 4.546       | 395           | 400         | 412          | rVB      | 239297         | 360098        | 7.04%           | 1.131%        |
| 3         | 4.751       | 427           | 435         | 449          | rVB3     | 28586          | 60211         | 1.18%           | 0.189%        |
| 4         | 5.104       | 485           | 495         | 501          | rBV2     | 104648         | 183352        | 3.58%           | 0.576%        |
| 5         | 5.381       | 535           | 542         | 546          | rBV      | 288718         | 394812        | 7.71%           | 1.240%        |
| 6         | 5.510       | 558           | 564         | 570          | rBV      | 929246         | 1235232       | 24.14%          | 3.879%        |
| 7         | 6.034       | 648           | 653         | 659          | rBV      | 41293          | 69811         | 1.36%           | 0.219%        |
| 8         | 6.098       | 659           | 664         | 669          | rVB4     | 46874          | 73824         | 1.44%           | 0.232%        |
| 9         | 6.410       | 708           | 717         | 722          | rBV      | 166230         | 404733        | 7.91%           | 1.271%        |
| 10        | 6.534       | 726           | 738         | 747          | rBV2     | 507246         | 1119091       | 21.87%          | 3.514%        |
| 11        | 6.704       | 760           | 767         | 775          | rBV4     | 77338          | 180187        | 3.52%           | 0.566%        |
| 12        | 6.792       | 775           | 782         | 786          | rBV2     | 54104          | 88245         | 1.72%           | 0.277%        |
| 13        | 6.892       | 793           | 799         | 804          | rBV      | 475197         | 638294        | 12.47%          | 2.004%        |
| 14        | 6.975       | 804           | 813         | 818          | rVV      | 2919284        | 5118003       | 100.00%         | 16.071%       |
| 15        | 7.045       | 821           | 825         | 828          | rVV2     | 75981          | 102387        | 2.00%           | 0.321%        |
| 16        | 7.092       | 828           | 833         | 837          | rVV3     | 80410          | 162164        | 3.17%           | 0.509%        |
| 17        | 7.134       | 837           | 840         | 844          | rVB3     | 73241          | 96140         | 1.88%           | 0.302%        |
| 18        | 7.292       | 863           | 867         | 872          | rVB4     | 59775          | 91616         | 1.79%           | 0.288%        |
| 19        | 7.457       | 882           | 895         | 899          | rBV      | 1208443        | 1837937       | 35.91%          | 5.771%        |
| 20        | 7.522       | 903           | 906         | 911          | rVB2     | 76119          | 107344        | 2.10%           | 0.337%        |
| 21        | 7.651       | 918           | 928         | 934          | rVB2     | 230199         | 623412        | 12.18%          | 1.958%        |
| 22        | 7.710       | 934           | 938         | 940          | rBV3     | 65134          | 86080         | 1.68%           | 0.270%        |
| 23        | 7.745       | 940           | 944         | 950          | rVB4     | 146202         | 249830        | 4.88%           | 0.784%        |
| 24        | 7.839       | 957           | 960         | 963          | rBV      | 57962          | 80007         | 1.56%           | 0.251%        |
| 25        | 7.916       | 968           | 973         | 977          | rBV2     | 112854         | 192071        | 3.75%           | 0.603%        |
| 26        | 8.069       | 992           | 999         | 1002         | rBV2     | 195794         | 333253        | 6.51%           | 1.046%        |
| 27        | 8.175       | 1011          | 1017        | 1022         | rVB2     | 801337         | 1276616       | 24.94%          | 4.009%        |
| 28        | 8.292       | 1033          | 1037        | 1048         | rVB4     | 133788         | 244463        | 4.78%           | 0.768%        |
| 29        | 8.398       | 1050          | 1055        | 1059         | rBV5     | 61983          | 147116        | 2.87%           | 0.462%        |
| 30        | 8.681       | 1099          | 1103        | 1109         | rVB      | 637821         | 853637        | 16.68%          | 2.680%        |
| 31        | 8.916       | 1140          | 1143        | 1145         | rBV3     | 104469         | 139735        | 2.73%           | 0.439%        |
| 32        | 9.098       | 1171          | 1174        | 1178         | rBV3     | 77767          | 124508        | 2.43%           | 0.391%        |
| 33        | 9.257       | 1196          | 1201        | 1205         | rBV      | 1939892        | 2515687       | 49.15%          | 7.899%        |
| 34        | 9.333       | 1211          | 1214        | 1215         | rBV3     | 85972          | 105874        | 2.07%           | 0.332%        |
| 35        | 9.357       | 1215          | 1218        | 1220         | rVV      | 144735         | 179573        | 3.51%           | 0.564%        |
| 36        | 9.651       | 1264          | 1268        | 1271         | rBV2     | 212136         | 319061        | 6.23%           | 1.002%        |

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Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

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-BF052025.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 9.933  | 1310 | 1316 | 1320 | rVB3 | 555772  | 1011901 | 19.77% | 3.177% |
| 38 | 10.733 | 1449 | 1452 | 1457 | rBV  | 620644  | 711035  | 13.89% | 2.233% |
| 39 | 10.857 | 1470 | 1473 | 1480 | rVB  | 938520  | 1383133 | 27.02% | 4.343% |
| 40 | 13.027 | 1838 | 1842 | 1850 | rVB  | 1461792 | 2183119 | 42.66% | 6.855% |
| 41 | 14.715 | 2125 | 2129 | 2134 | rBV2 | 1613818 | 2885442 | 56.38% | 9.060% |
| 42 | 16.309 | 2395 | 2400 | 2417 | rVB  | 678888  | 1885096 | 36.83% | 5.919% |
| 43 | 16.762 | 2471 | 2477 | 2489 | rVB  | 888335  | 1834046 | 35.84% | 5.759% |

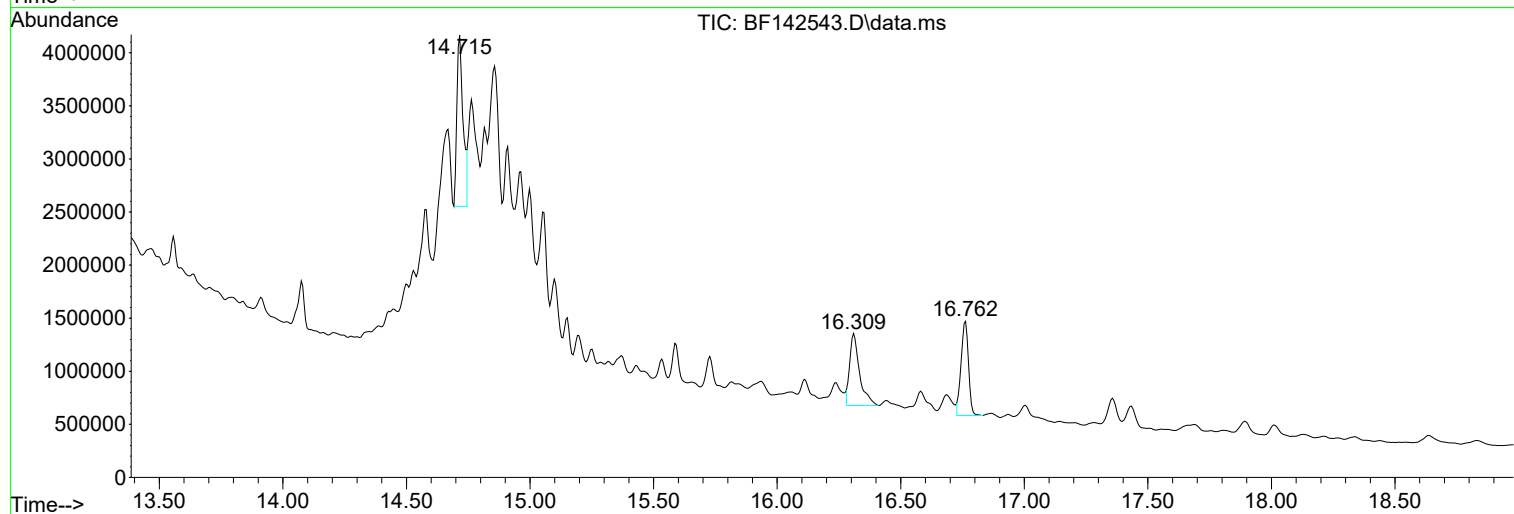
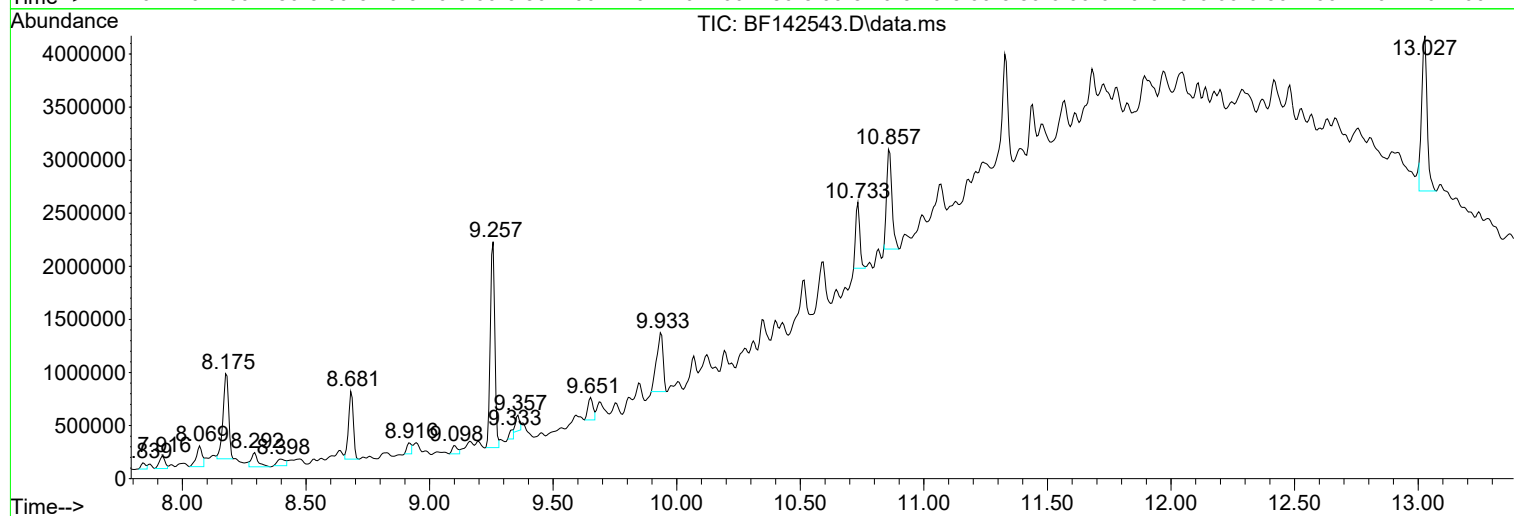
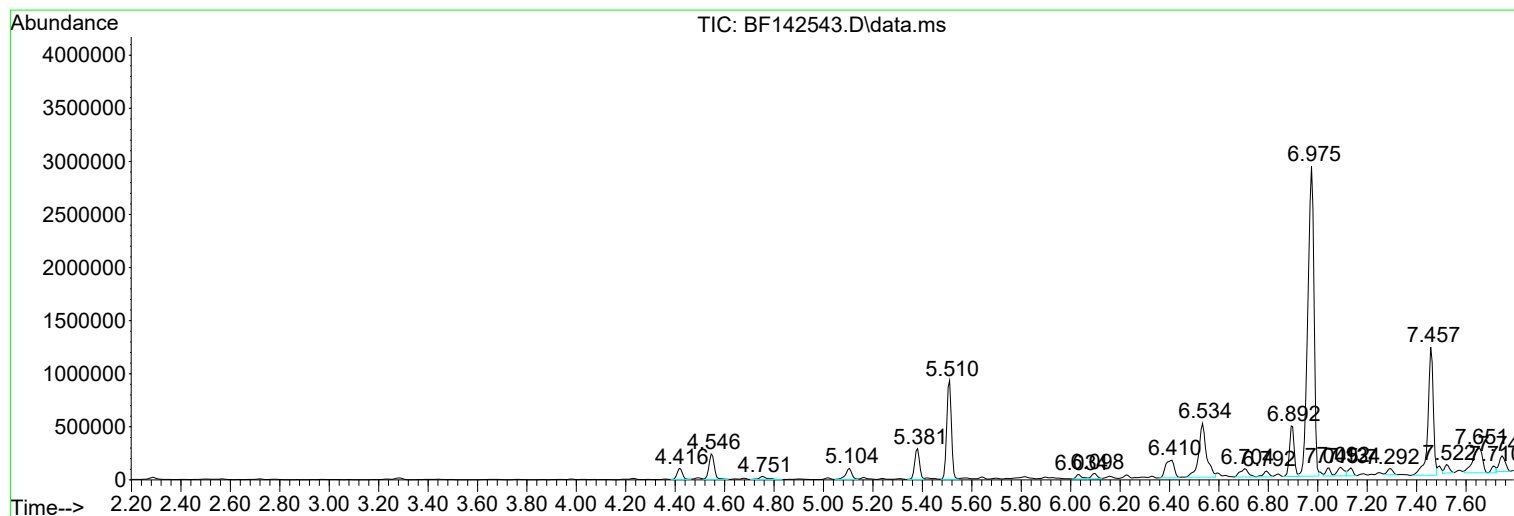
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
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Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.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\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

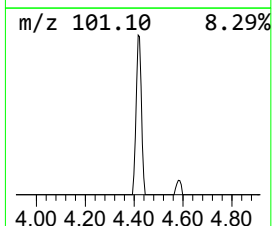
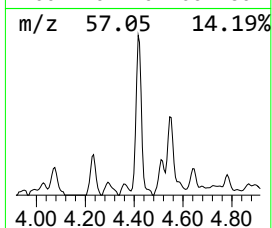
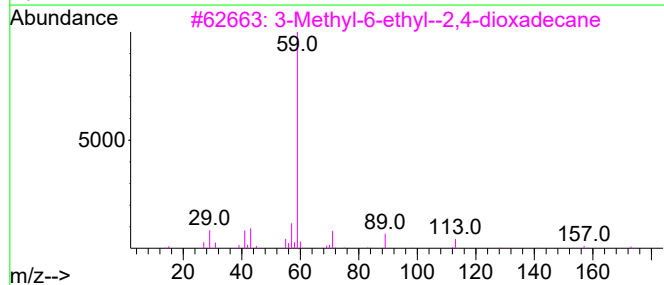
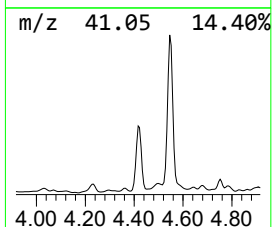
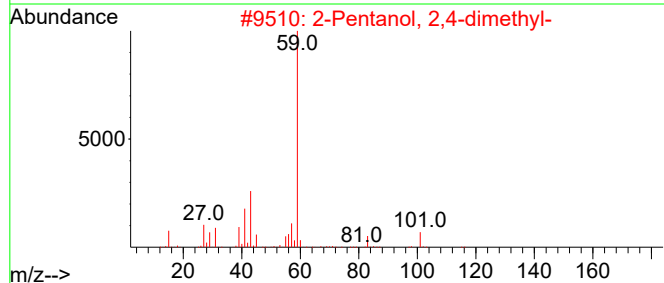
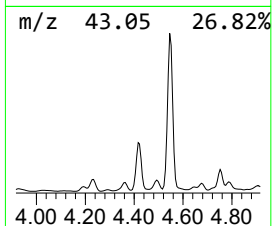
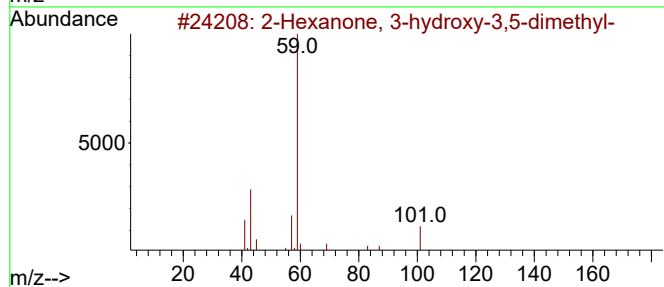
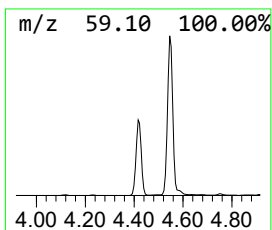
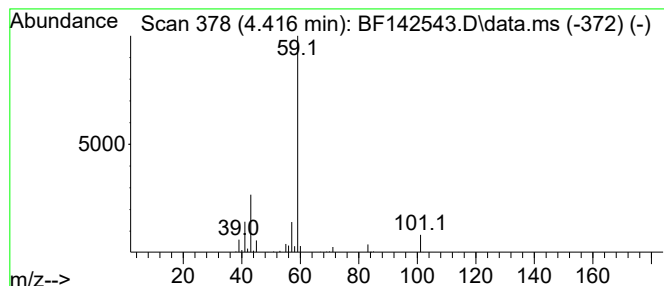
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 2-Hexanone, 3-hydroxy-3,5-d... Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 4.416 | 4.97 ng | 158720 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2-Hexanone, 3-hydroxy-3,5-dimethyl- | 144 | C8H16O2      | 006321-14-8  | 78      |      |      |
| 2    | 2-Pentanol, 2,4-dimethyl-           | 116 | C7H16O       | 000625-06-9  | 56      |      |      |
| 3    | 3-Methyl-6-ethyl--2,4-dioxadecane   | 188 | C11H24O2     | 1000334-83-2 | 56      |      |      |
| 4    | 4-Pentene-2-ol, 2-methyl            | 100 | C6H12O       | 000624-97-5  | 42      |      |      |
| 5    | Propane, 1-ethoxy-2-methyl-         | 102 | C6H14O       | 000627-02-1  | 40      |      |      |



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Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

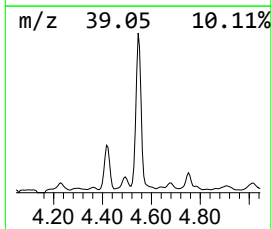
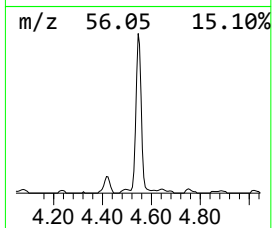
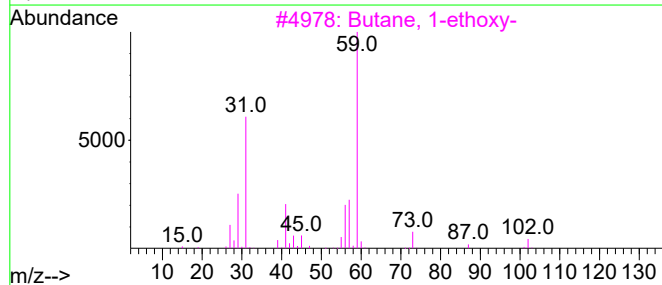
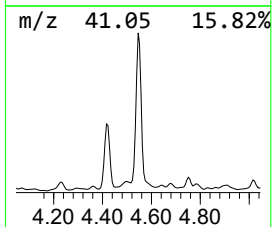
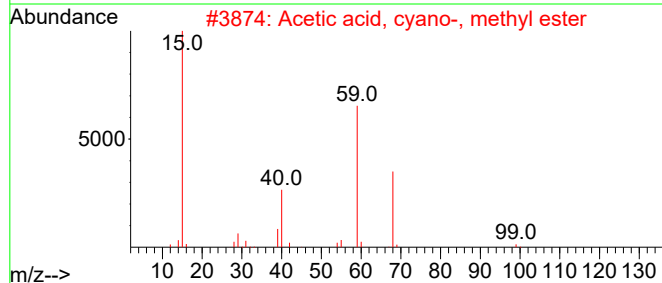
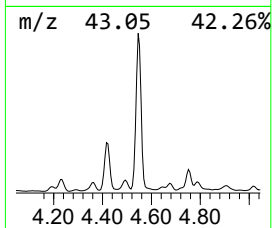
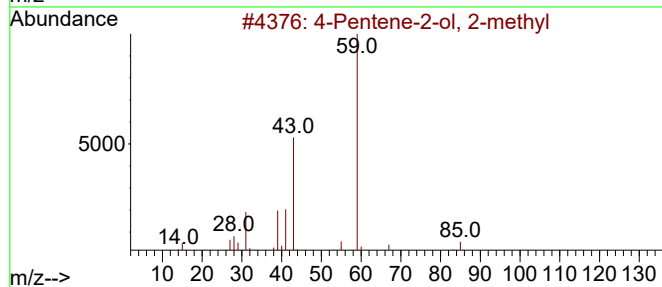
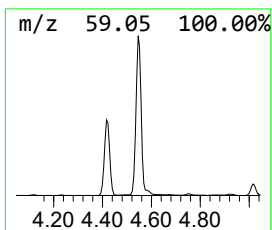
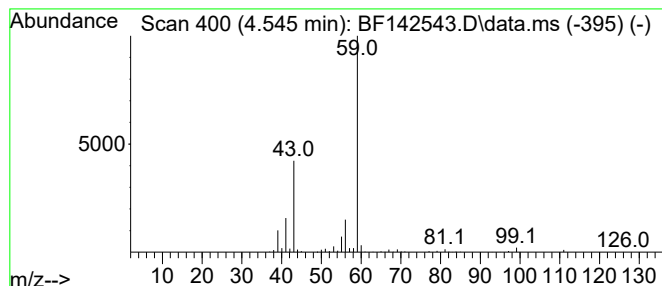
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 4-Pentene-2-ol, 2-methyl Concentration Rank 8

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 4.545 | 11.28 ng | 360098 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 4-Pentene-2-ol, 2-methyl            | 100 | C6H12O  | 000624-97-5 | 50   |
| 2    |    |   | Acetic acid, cyano-, methyl ester   | 99  | C4H5NO2 | 000105-34-0 | 45   |
| 3    |    |   | Butane, 1-ethoxy-                   | 102 | C6H14O  | 000628-81-9 | 39   |
| 4    |    |   | Propanoic acid, 2-nitro-, methyl... | 133 | C4H7NO4 | 006118-50-9 | 39   |
| 5    |    |   | Oxirane, tetramethyl-               | 100 | C6H12O  | 005076-20-0 | 9    |



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ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
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LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

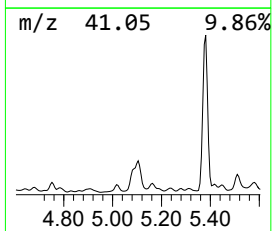
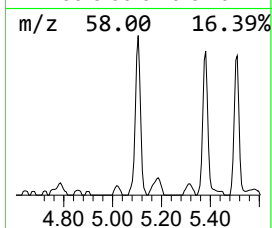
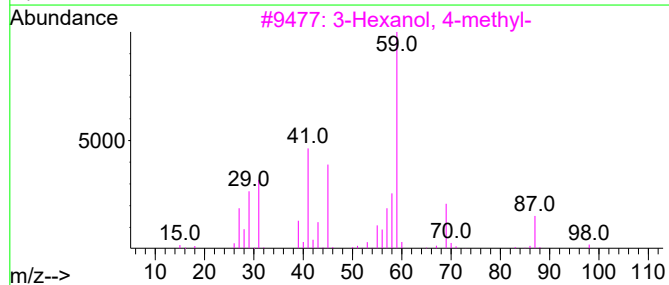
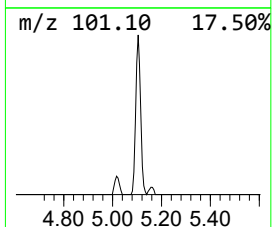
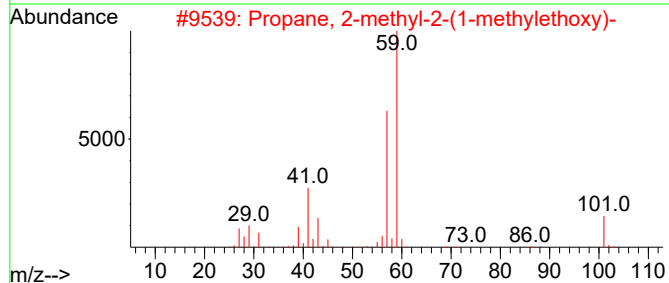
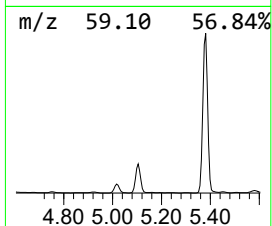
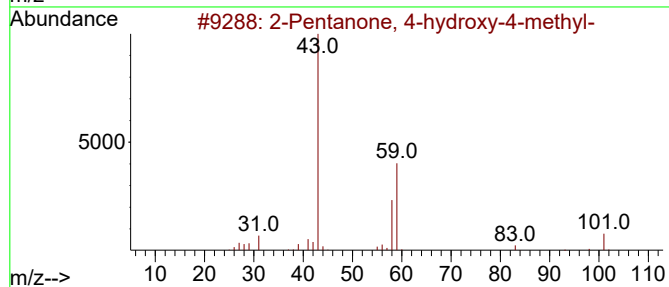
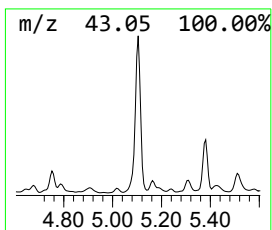
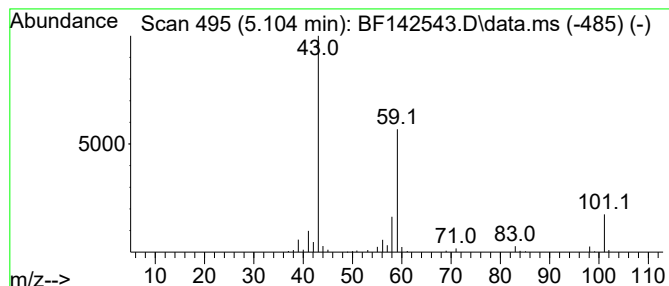
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 5.104 | 5.75 ng | 183352 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 45   |
| 2    |    |   | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O   | 017348-59-3 | 42   |
| 3    |    |   | 3-Hexanol, 4-methyl-                | 116 | C7H16O   | 000615-29-2 | 33   |
| 4    |    |   | 2-Hexanol, 2-methyl-                | 116 | C7H16O   | 000625-23-0 | 28   |
| 5    |    |   | 1,8-Nonanediol, 8-methyl-           | 174 | C10H22O2 | 054725-73-4 | 25   |



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ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
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LAW-25-0078

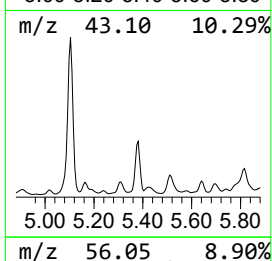
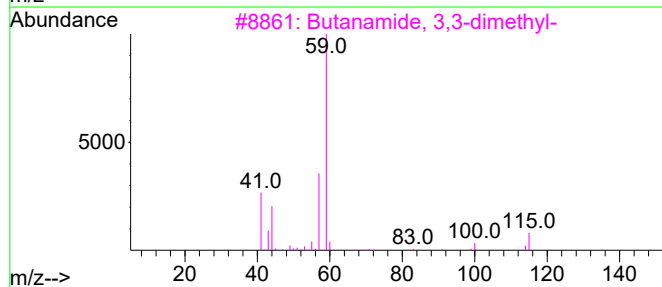
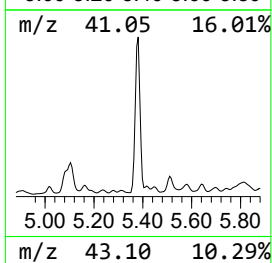
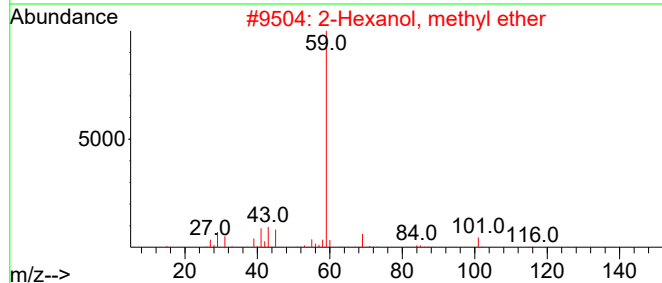
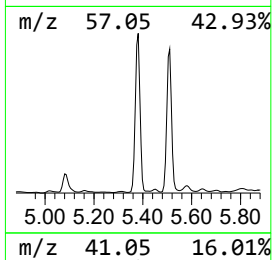
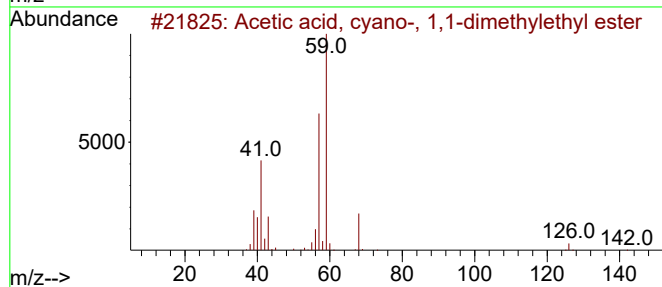
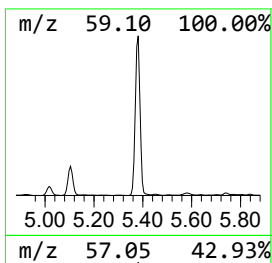
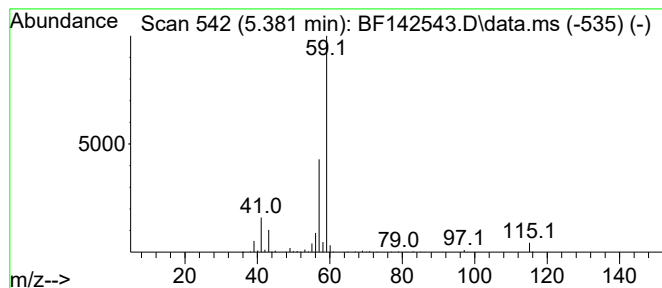
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Acetic acid, cyano-, 1,1-di... Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 5.381 | 12.37 ng | 394812 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9  | 64   |
| 2    |    |   | 2-Hexanol, methyl ether             | 116 | C7H16O   | 1000333-92-0 | 9    |
| 3    |    |   | Butanamide, 3,3-dimethyl-           | 115 | C6H13NO  | 000926-04-5  | 9    |
| 4    |    |   | Propane, 2-ethoxy-2-methyl-         | 102 | C6H14O   | 000637-92-3  | 9    |
| 5    |    |   | 2-Propanol, 2-methyl-               | 74  | C4H10O   | 000075-65-0  | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

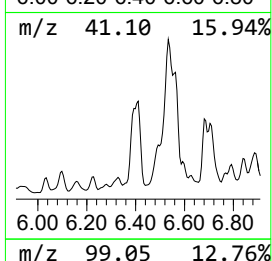
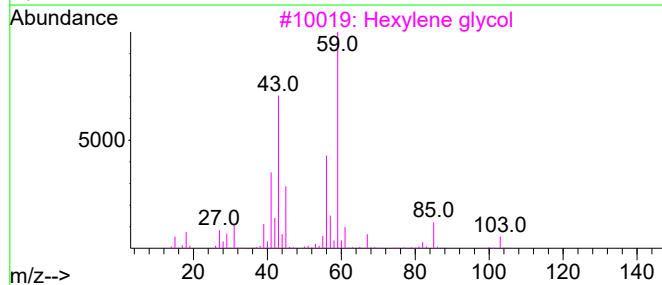
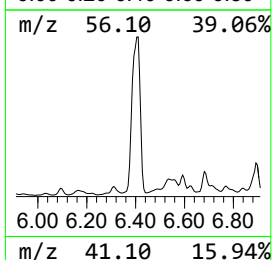
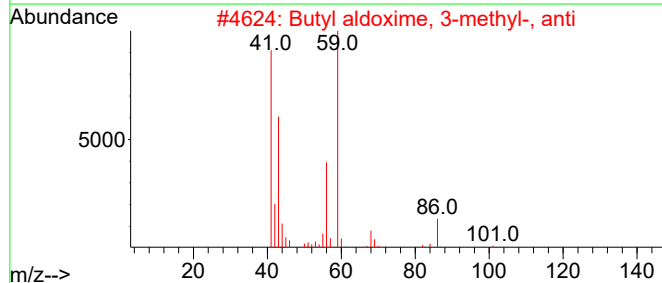
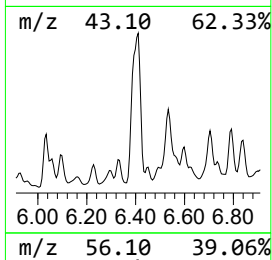
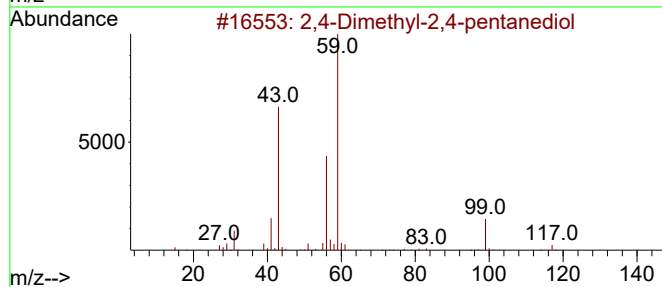
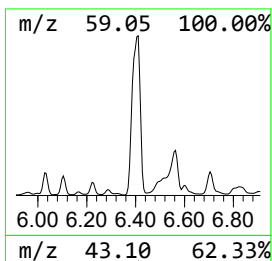
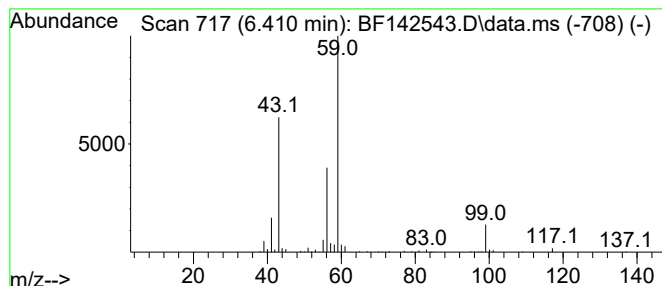
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 2,4-Dimethyl-2,4-pentanediol Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 6.410 | 12.68 ng | 404733 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2,4-Dimethyl-2,4-pentanediol        | 132 | C7H16O2 | 024892-49-7 | 83   |
| 2    |    |   | Butyl aldoxime, 3-methyl-, anti     | 101 | C5H11NO | 005775-74-6 | 72   |
| 3    |    |   | Hexylene glycol                     | 118 | C6H14O2 | 000107-41-5 | 56   |
| 4    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 38   |
| 5    |    |   | Oxetane, 2,2,4-trimethyl-           | 100 | C6H12O  | 023120-44-7 | 36   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

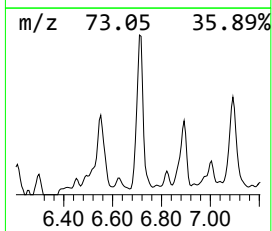
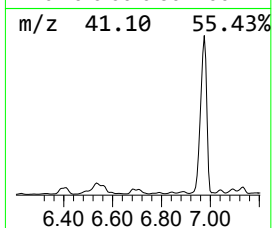
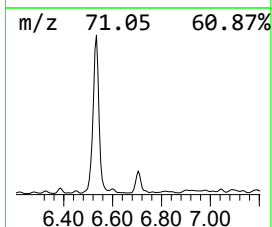
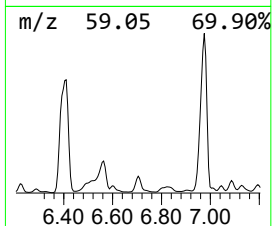
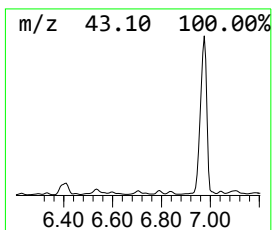
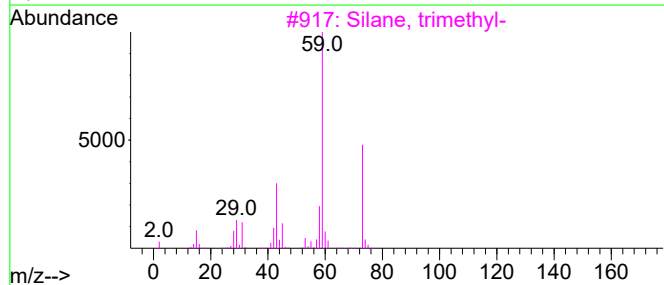
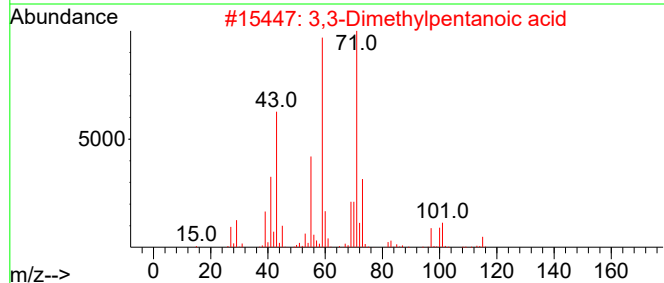
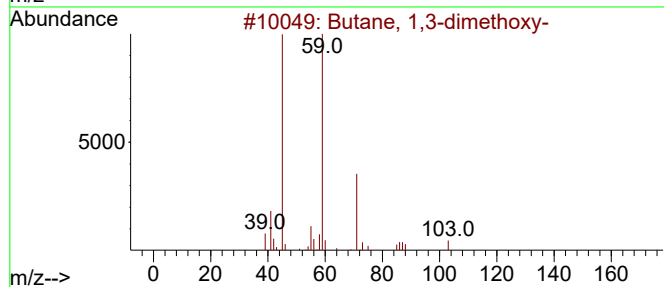
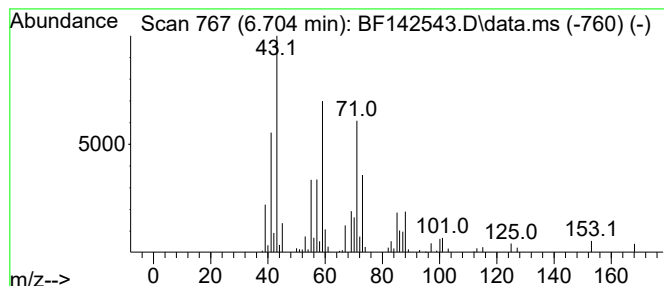
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Butane, 1,3-dimethoxy- Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.704 | 5.65 ng | 180187 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 1,3-dimethoxy-         | 118 | C6H14O2 | 010143-66-5 | 59   |
| 2    |    |   | 3,3-Dimethylpentanoic acid     | 130 | C7H14O2 | 003177-74-0 | 50   |
| 3    |    |   | Silane, trimethyl-             | 74  | C3H10Si | 000993-07-7 | 27   |
| 4    |    |   | 1-Butene, 3-butoxy-2-methyl-   | 142 | C9H18O  | 056585-28-5 | 22   |
| 5    |    |   | 1-Propene, 1-methoxy-2-methyl- | 86  | C5H10O  | 017574-84-4 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

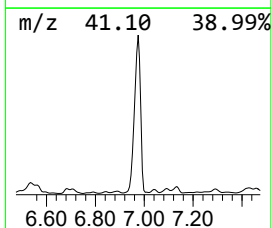
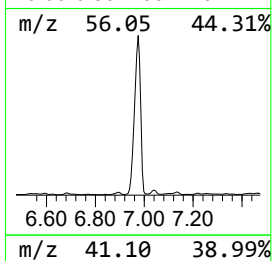
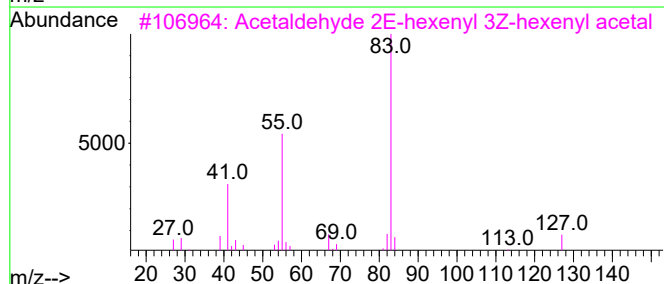
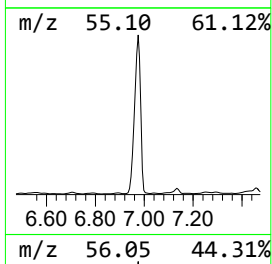
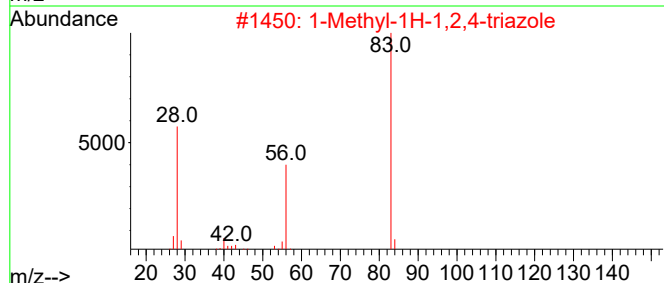
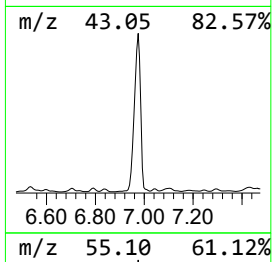
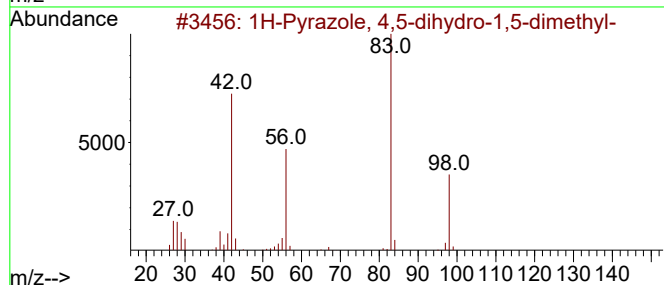
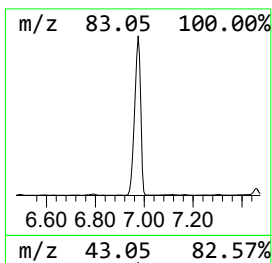
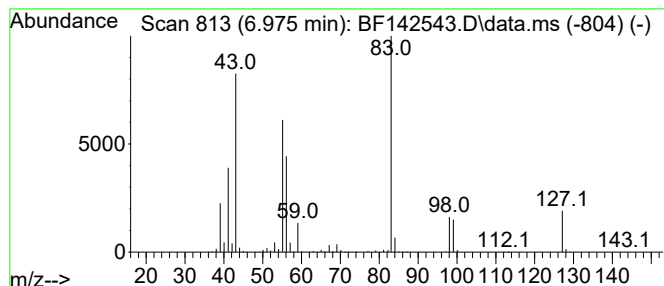
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 1H-Pyrazole, 4,5-dihydro-1,... Concentration Rank 1

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 6.975 | 160.37 ng | 5118000 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|------|-------------------------------------|-----|----------|--------------|------|
| 1    |      | 1H-Pyrazole, 4,5-dihydro-1,5-dim... | 98  | C5H10N2  | 005775-96-2  | 50   |
| 2    |      | 1-Methyl-1H-1,2,4-triazole          | 83  | C3H5N3   | 006086-21-1  | 47   |
| 3    |      | Acetaldehyde 2E-hexenyl 3Z-hexen... | 226 | C14H26O2 | 1000431-45-5 | 38   |
| 4    |      | 1-Heptanol, 2,4-dimethyl-,          | 144 | C9H20O   | 098982-97-9  | 37   |
| 5    |      | 2-Pentene, 4,4-dimethyl-            | 98  | C7H14    | 026232-98-4  | 37   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

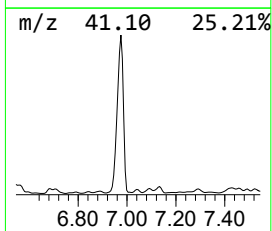
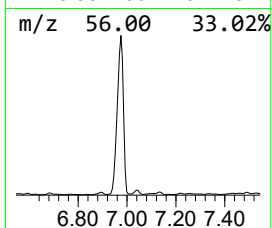
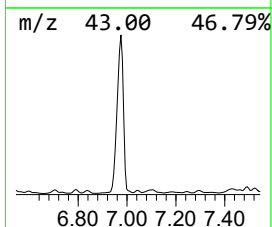
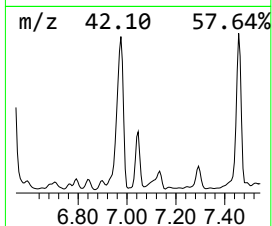
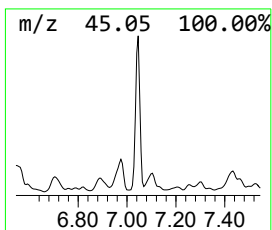
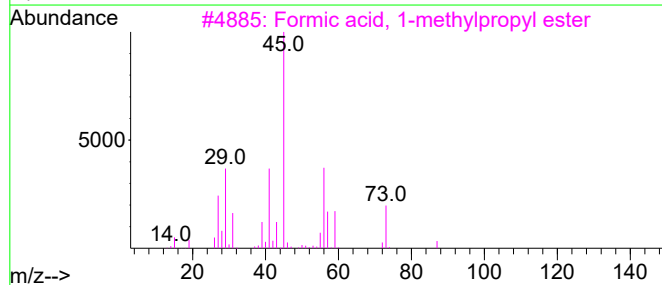
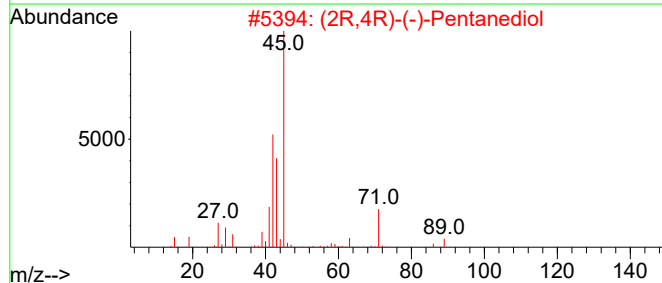
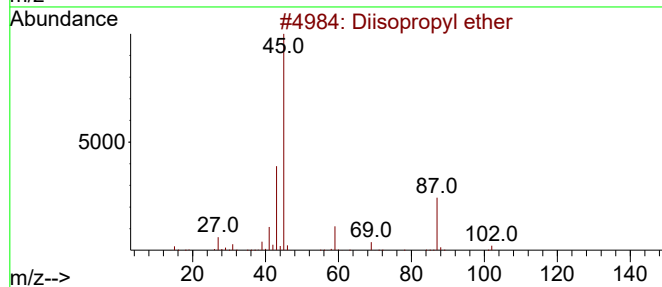
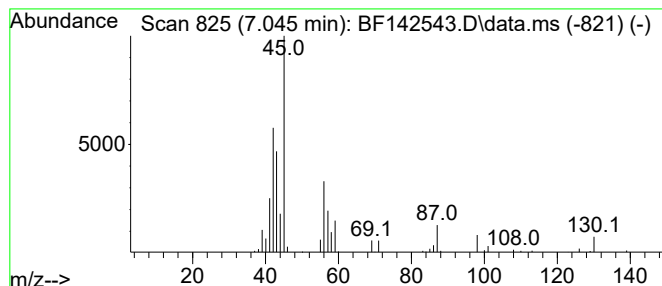
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 unknown7.045 Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 7.045 | 3.21 ng | 102387 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Diisopropyl ether                   | 102 | C6H14O   | 000108-20-3 | 43   |
| 2    |    |   | (2R,4R)-(-)-Pentanediol             | 104 | C5H12O2  | 042075-32-1 | 40   |
| 3    |    |   | Formic acid, 1-methylpropyl ester   | 102 | C5H10O2  | 000589-40-2 | 38   |
| 4    |    |   | 1,4-Pentanediol                     | 104 | C5H12O2  | 000626-95-9 | 38   |
| 5    |    |   | Ethanol, 2-[2-(2-butoxyethoxy)et... | 206 | C10H22O4 | 000143-22-6 | 37   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

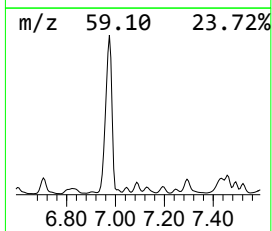
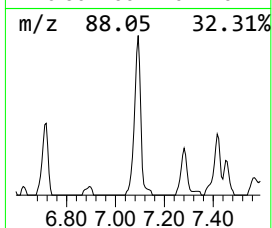
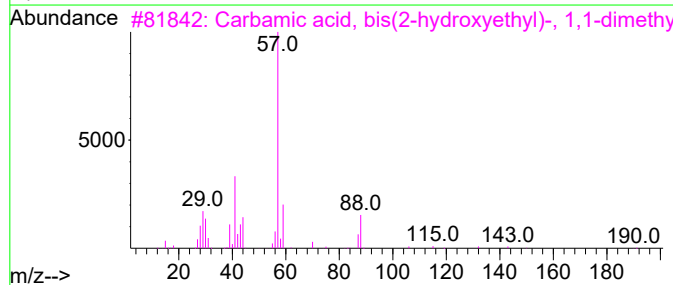
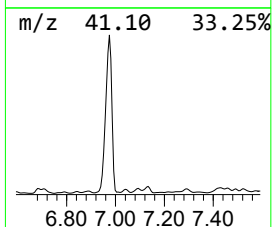
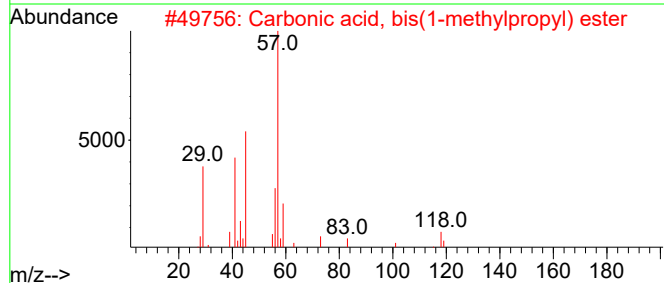
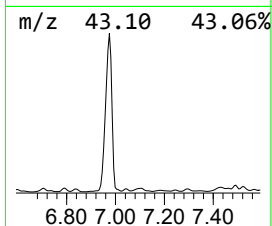
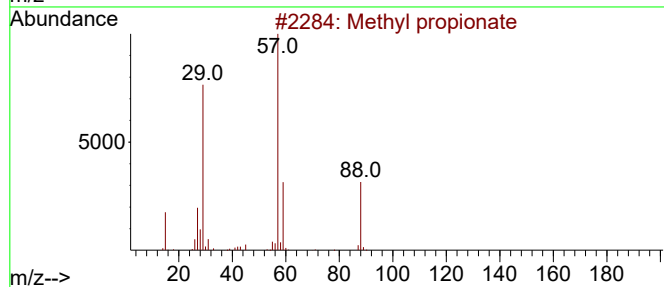
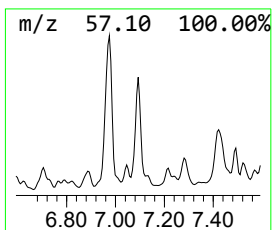
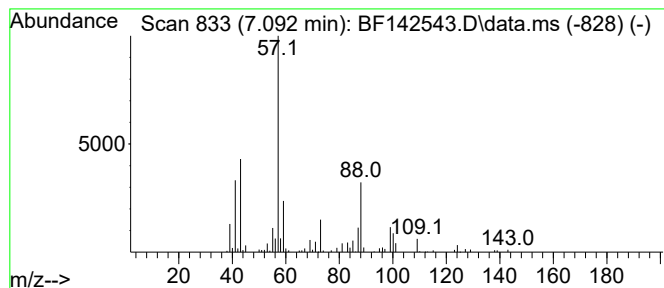
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 unknown7.092 Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 7.092 | 5.08 ng | 162164 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Methyl propionate                   | 88  | C4H8O2    | 000554-12-1  | 46   |
| 2    |    |   | Carbonic acid, bis(1-methylpropy... | 174 | C9H18O3   | 000623-63-2  | 35   |
| 3    |    |   | Carbamic acid, bis(2-hydroxyethy... | 205 | C9H19NO4  | 103898-11-9  | 25   |
| 4    |    |   | Hydrazinecarboxylic acid, 1,1-di... | 132 | C5H12N2O2 | 000870-46-2  | 16   |
| 5    |    |   | 1,3-Dioxolane, 2,2-dimethyl-4-(t... | 188 | C10H20O3  | 1000381-76-2 | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

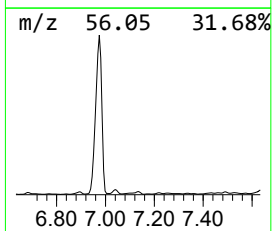
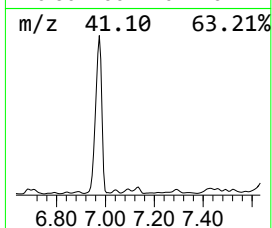
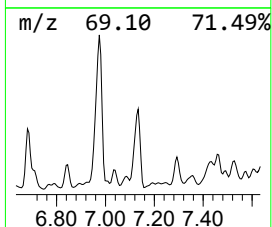
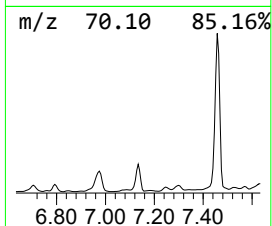
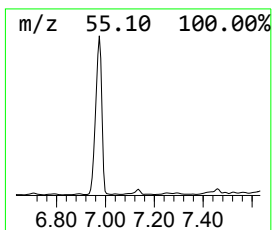
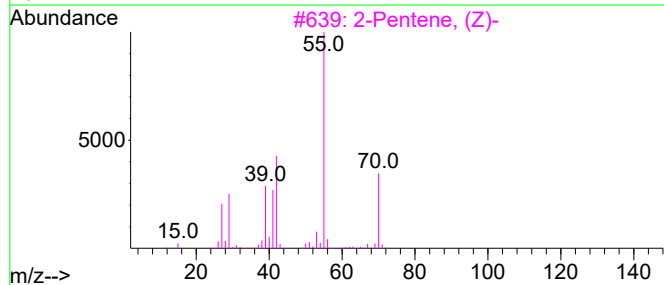
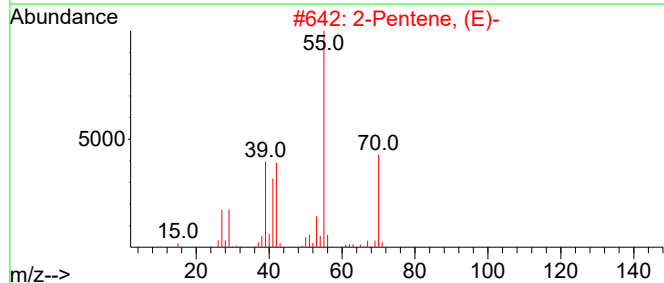
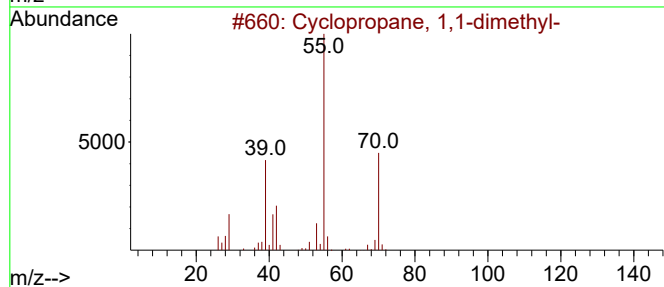
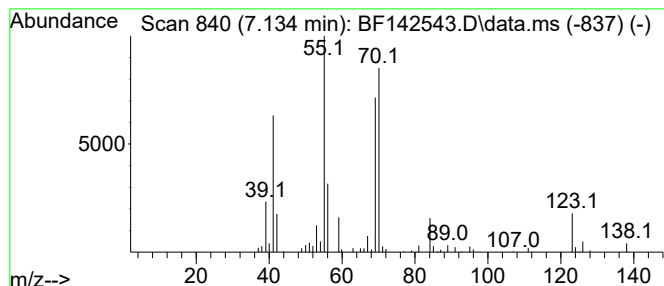
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Cyclopropane, 1,1-dimethyl- Concentration Rank 20

| R.T.  | EstConc | Area  | Relative to ISTD       | R.T.  |
|-------|---------|-------|------------------------|-------|
| 7.134 | 3.01 ng | 96140 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                        | MW | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|----|---------|-------------|------|
| 1    |    |   | Cyclopropane, 1,1-dimethyl-         | 70 | C5H10   | 001630-94-0 | 50   |
| 2    |    |   | 2-Pentene, (E)-                     | 70 | C5H10   | 000646-04-8 | 50   |
| 3    |    |   | 2-Pentene, (Z)-                     | 70 | C5H10   | 000627-20-3 | 43   |
| 4    |    |   | 2-Pentene                           | 70 | C5H10   | 000109-68-2 | 43   |
| 5    |    |   | Cyclopropane, 1,2-dimethyl-, trans- | 70 | C5H10   | 002402-06-4 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

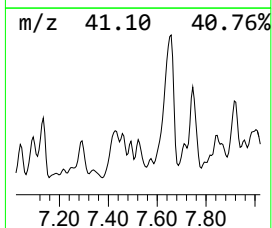
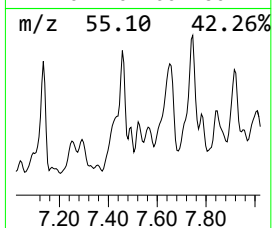
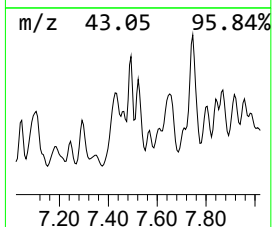
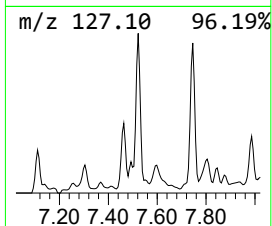
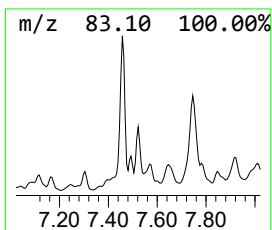
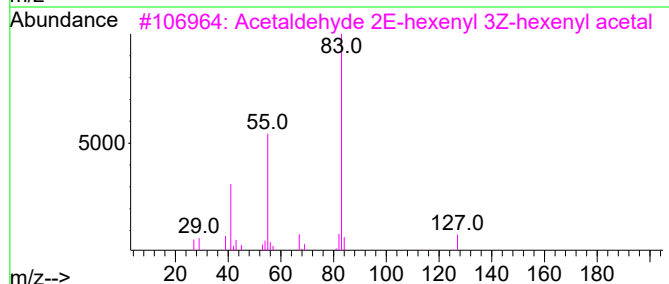
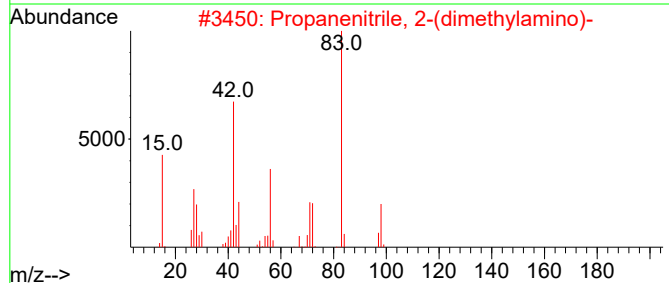
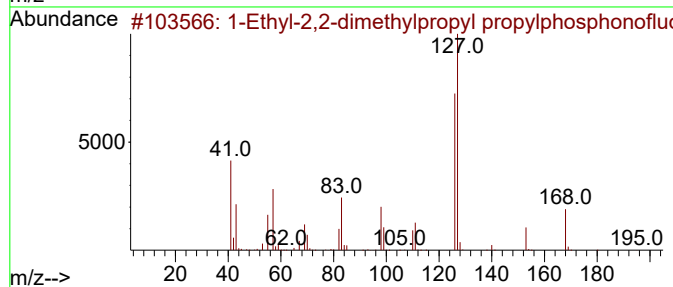
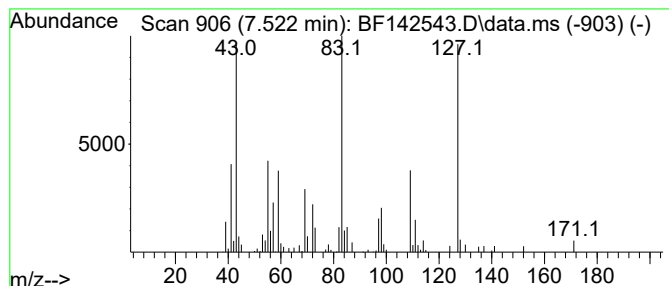
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 unknown7.522 Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 7.522 | 3.36 ng | 107344 | 1,4-Dichlorobenzene-d4 | 6.898 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 1-Ethyl-2,2-dimethylpropyl propy... | 224 | C10H22F02P | 1010298-29-2 | 38   |
| 2    |    |   | Propanenitrile, 2-(dimethylamino)-  | 98  | C5H10N2    | 005350-67-4  | 35   |
| 3    |    |   | Acetaldehyde 2E-hexenyl 3Z-hexen... | 226 | C14H26O2   | 1000431-45-5 | 27   |
| 4    |    |   | Isopropyl tiglate                   | 142 | C8H14O2    | 001733-25-1  | 27   |
| 5    |    |   | 1-Methyl-3-nitropyrazole            | 127 | C4H5N3O2   | 054210-32-1  | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

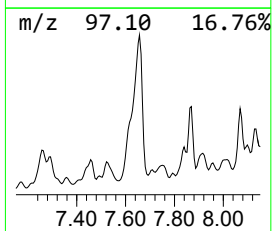
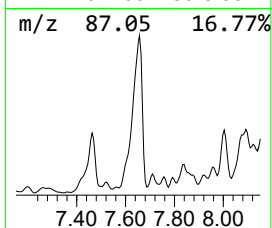
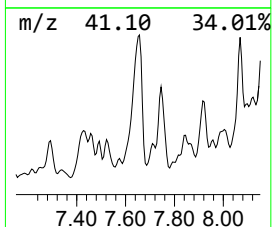
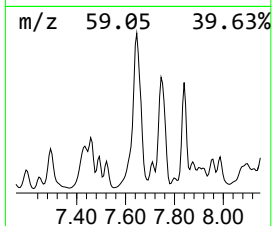
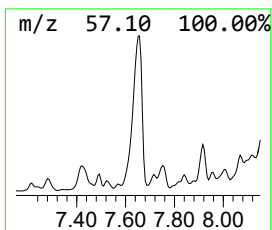
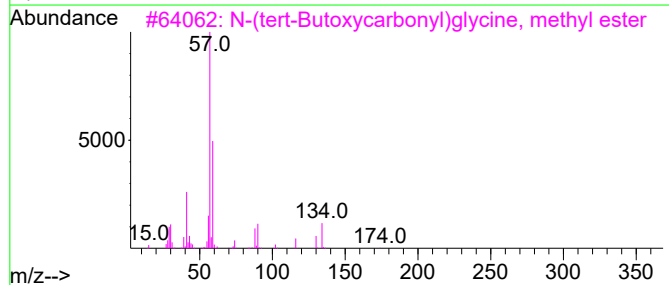
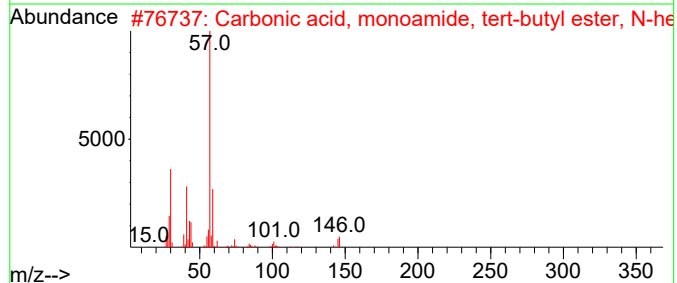
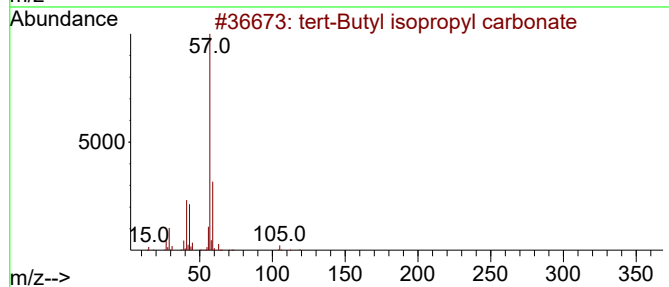
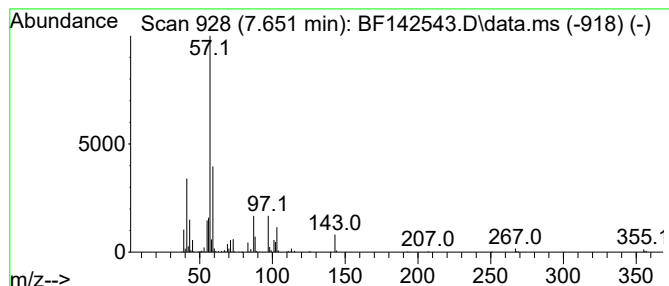
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 unknown7.651 Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 7.651 | 9.77 ng | 623412 | Naphthalene-d8   | 8.181 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | tert-Butyl isopropyl carbonate      | 160 | C8H16O3   | 030221-21-7  | 38   |
| 2    |    |   | Carbonic acid, monoamide, tert-b... | 201 | C11H23NO2 | 1000420-59-4 | 37   |
| 3    |    |   | N-(tert-Butoxycarbonyl)glycine, ... | 189 | C8H15NO4  | 1000333-18-9 | 37   |
| 4    |    |   | Propane, 1-(1,1-dimethylethoxy)-... | 130 | C8H18O    | 033021-02-2  | 32   |
| 5    |    |   | 3-tert-Butoxycarbonylamino-butyr... | 203 | C9H17NO4  | 1000317-53-1 | 28   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
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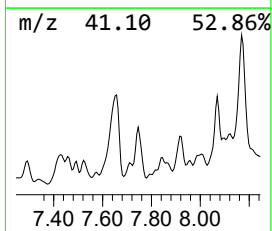
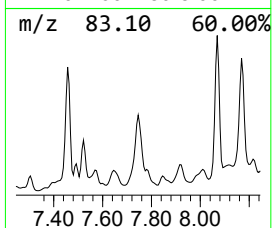
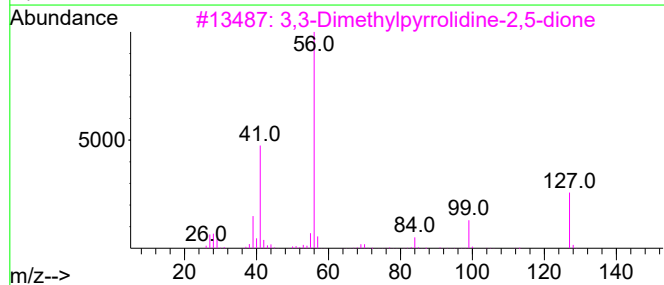
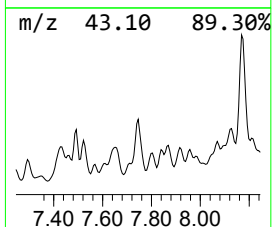
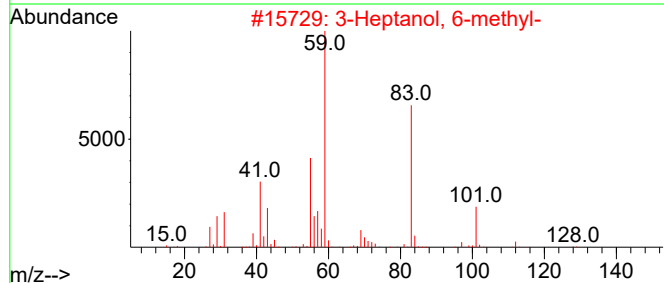
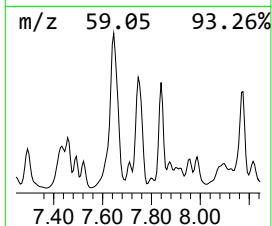
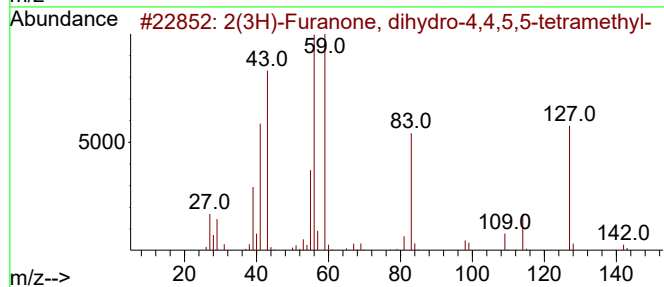
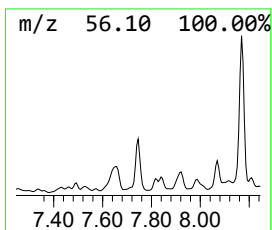
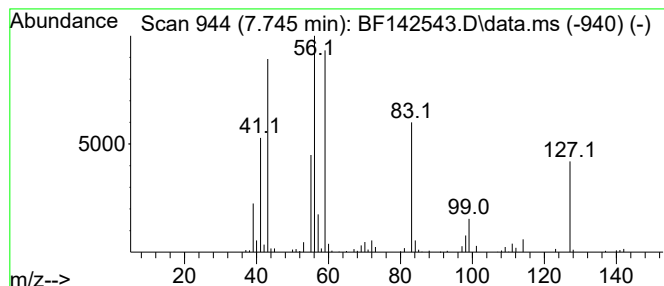
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Peak Number 13 2(3H)-Furanone, dihydro-4,4... Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 7.745 | 3.91 ng | 249830 | Naphthalene-d8   | 8.181 |

| Hit# | of 5                                  | Tentative ID | MW      | MolForm | CAS#        | Qual |
|------|---------------------------------------|--------------|---------|---------|-------------|------|
| 1    | 2(3H)-Furanone, dihydro-4,4,5,5-...   | 142          | C8H14O2 |         | 016466-24-3 | 78   |
| 2    | 3-Heptanol, 6-methyl-                 | 130          | C8H18O  |         | 018720-66-6 | 38   |
| 3    | 3,3-Dimethylpyrrolidine-2,5-dione     | 127          | C6H9NO2 |         | 003437-29-4 | 38   |
| 4    | 2,2,5,5-Tetramethyltetrahydro-3-...   | 142          | C8H14O2 |         | 005455-94-7 | 37   |
| 5    | 1H-1,2,4-Triazole-3-carboxaldehyde... | 111          | C4H5N3O |         | 056804-98-9 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

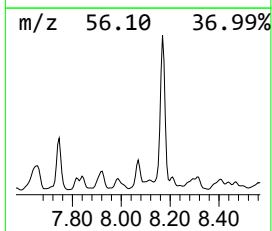
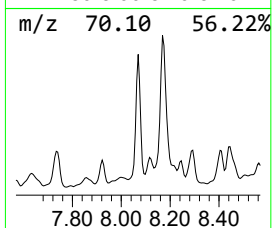
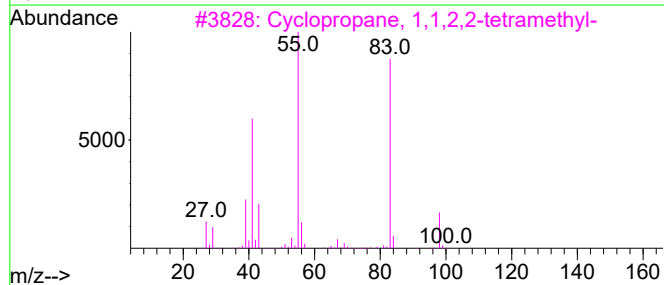
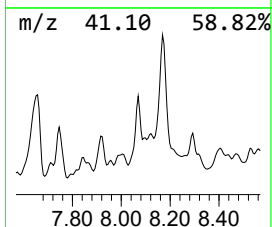
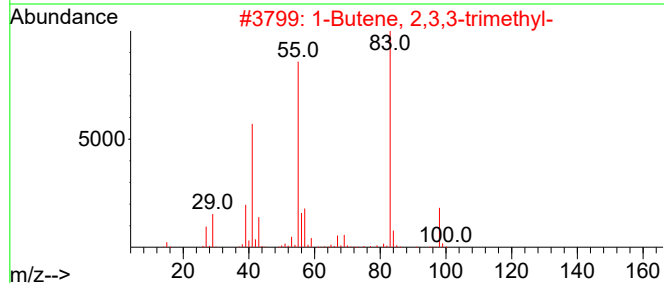
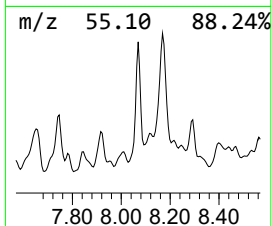
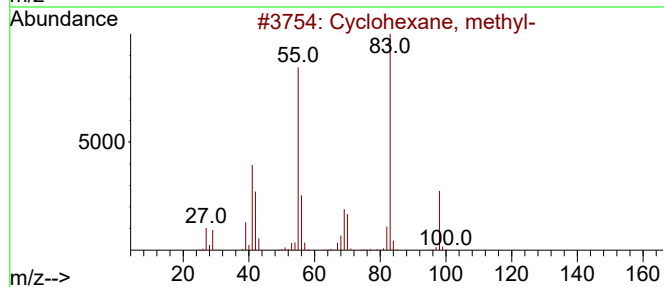
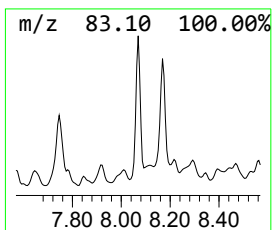
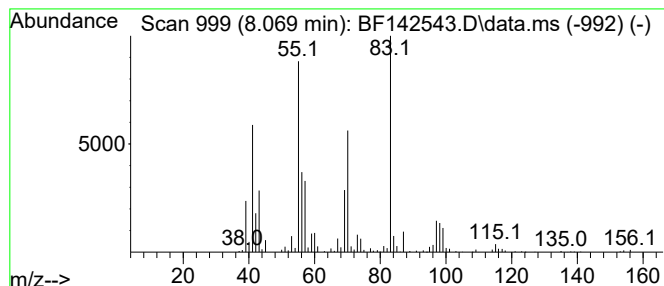
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 Cyclohexane, methyl- Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 8.069 | 5.22 ng | 333253 | Naphthalene-d8   | 8.181 |

| Hit# | of 5 | Tentative ID                       | MW | MolForm | CAS#        | Qual |
|------|------|------------------------------------|----|---------|-------------|------|
| 1    |      | Cyclohexane, methyl-               | 98 | C7H14   | 000108-87-2 | 58   |
| 2    |      | 1-Butene, 2,3,3-trimethyl-         | 98 | C7H14   | 000594-56-9 | 49   |
| 3    |      | Cyclopropane, 1,1,2,2-tetramethyl- | 98 | C7H14   | 004127-47-3 | 49   |
| 4    |      | 2-Pentene, 4,4-dimethyl-, (Z)-     | 98 | C7H14   | 000762-63-0 | 46   |
| 5    |      | 1,4-Pentadien-3-ol                 | 84 | C5H8O   | 000922-65-6 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

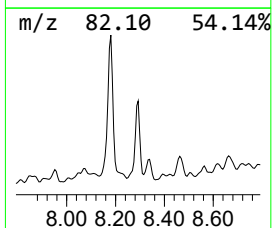
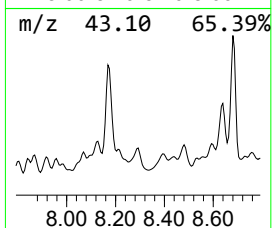
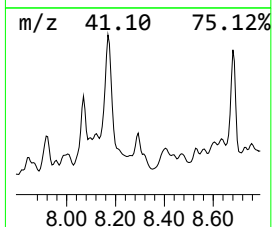
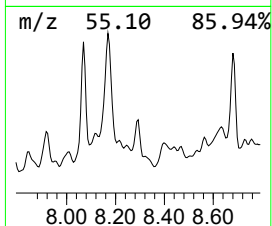
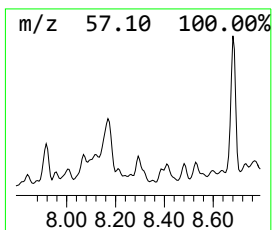
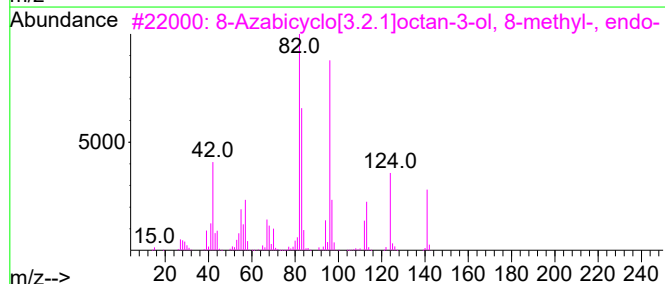
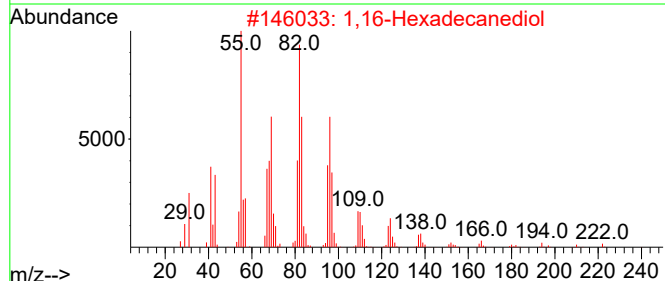
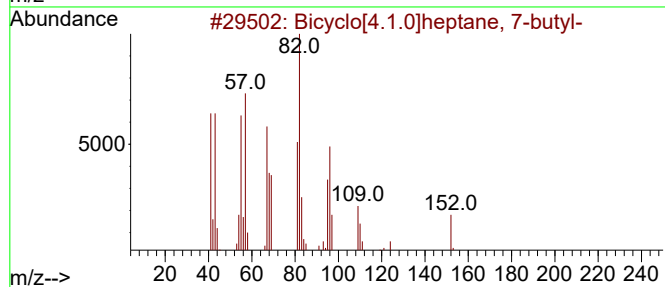
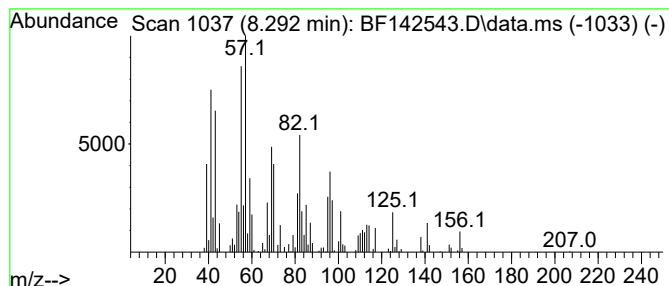
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 unknown8.292 Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 8.292 | 3.83 ng | 244463 | Naphthalene-d8   | 8.181 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Bicyclo[4.1.0]heptane, 7-butyl-     | 152 | C11H20   | 018645-10-8  | 41   |
| 2    |    |   | 1,16-Hexadecanediol                 | 258 | C16H34O2 | 007735-42-4  | 35   |
| 3    |    |   | 8-Azabicyclo[3.2.1]octan-3-ol, 8... | 141 | C8H15NO  | 000120-29-6  | 30   |
| 4    |    |   | Dihydroterrein                      | 156 | C8H12O3  | 1350310-16-5 | 27   |
| 5    |    |   | 3-Cyclopentene-1-acetaldehyde, 2... | 124 | C7H8O2   | 051905-39-6  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

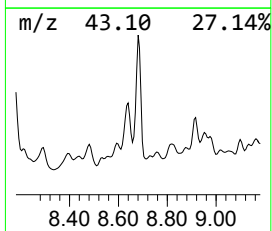
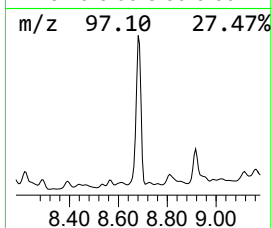
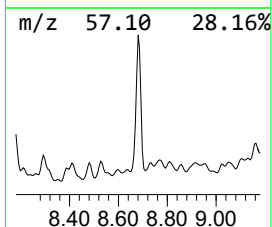
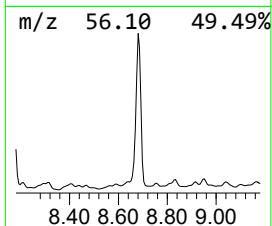
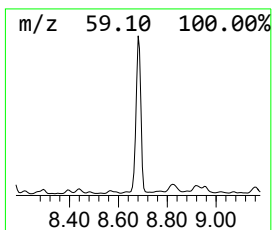
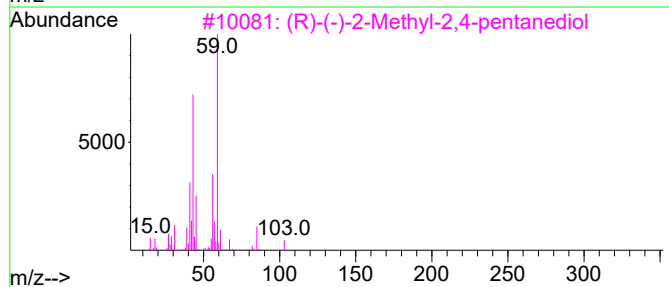
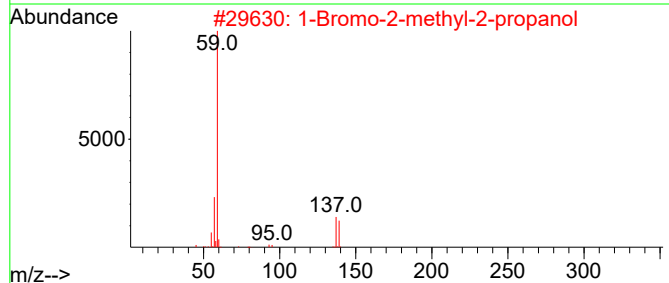
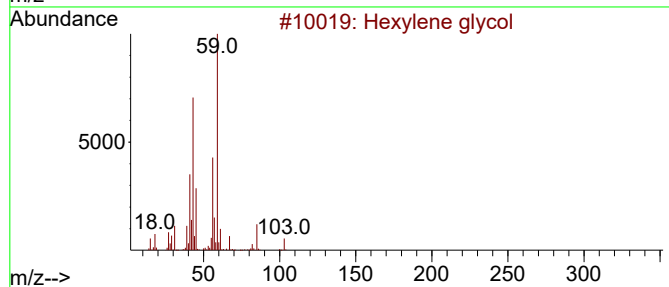
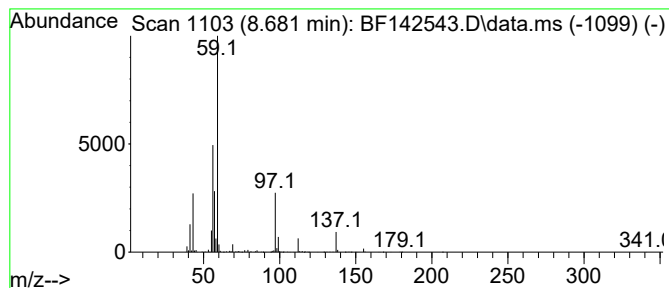
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 unknown8.681 Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 8.681 | 13.37 ng | 853637 | Naphthalene-d8   | 8.181 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexylene glycol                     | 118 | C6H14O2 | 000107-41-5 | 42   |
| 2    |    |   | 1-Bromo-2-methyl-2-propanol         | 152 | C4H9BrO | 038254-49-8 | 37   |
| 3    |    |   | (R)-(-)-2-Methyl-2,4-pentanediol    | 118 | C6H14O2 | 099210-90-9 | 36   |
| 4    |    |   | Propanoic acid, 2-hydroxy-2-meth... | 118 | C5H10O3 | 002110-78-3 | 32   |
| 5    |    |   | Hexanamide                          | 115 | C6H13NO | 000628-02-4 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

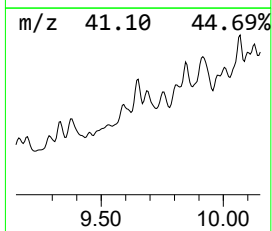
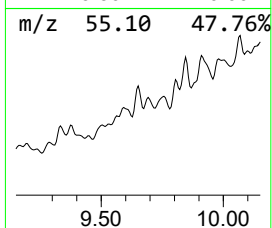
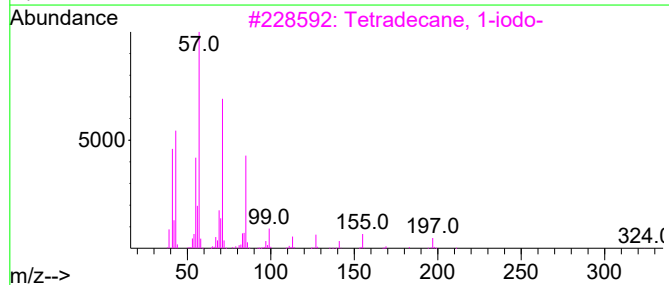
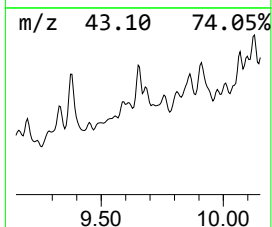
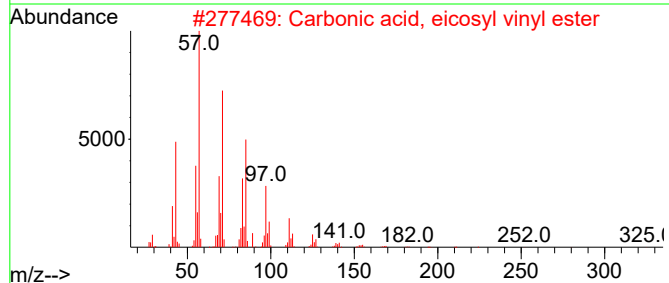
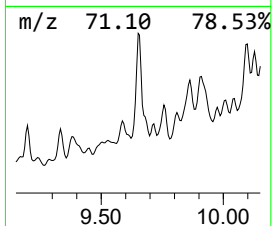
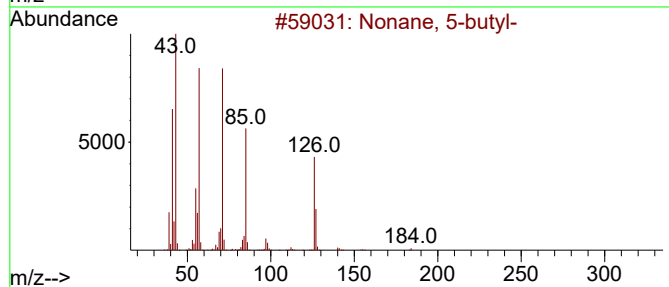
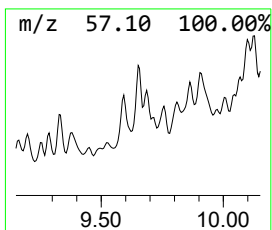
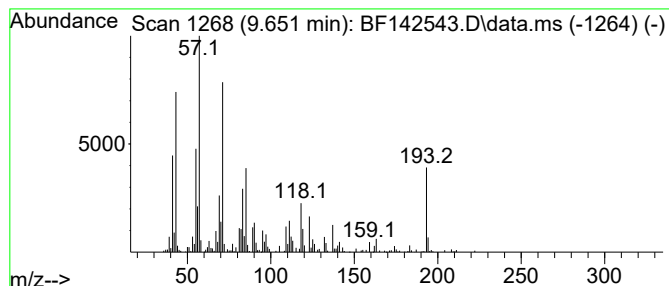
Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.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 Nonane, 5-butyl- Concentration Rank 10

| R.T.      | EstConc                            | Area   | Relative to ISTD |          | R.T.         |      |
|-----------|------------------------------------|--------|------------------|----------|--------------|------|
| 9.651     | 6.31 ng                            | 319061 | Acenaphthene-d10 |          | 9.939        |      |
| Hit# of 5 | Tentative ID                       |        | MW               | MolForm  | CAS#         | Qual |
| 1         | Nonane, 5-butyl-                   |        | 184              | C13H28   | 017312-63-9  | 50   |
| 2         | Carbonic acid, eicosyl vinyl ester |        | 368              | C23H44O3 | 1000382-54-3 | 46   |
| 3         | Tetradecane, 1-iodo-               |        | 324              | C14H29I  | 019218-94-1  | 43   |
| 4         | Dodecane, 4-methyl-                |        | 184              | C13H28   | 006117-97-1  | 43   |
| 5         | Tetradecane                        |        | 198              | C14H30   | 000629-59-4  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

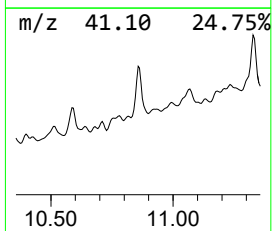
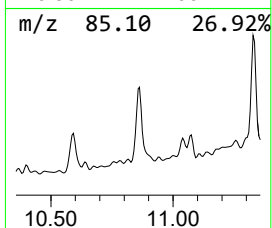
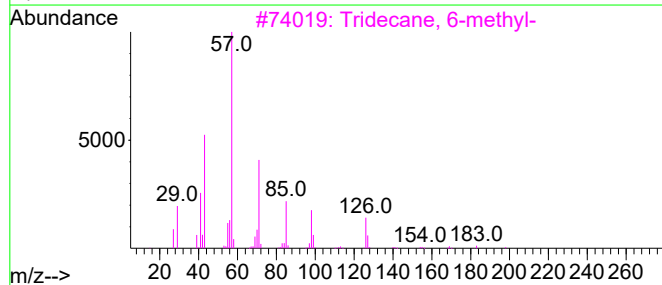
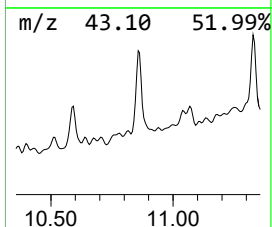
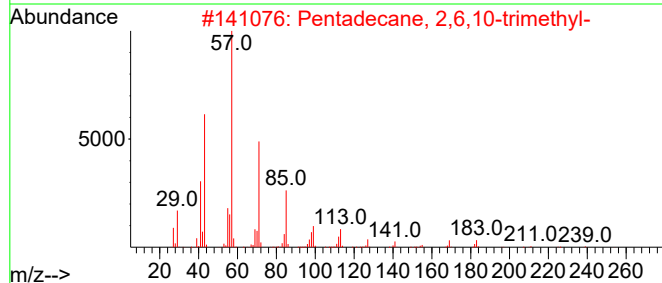
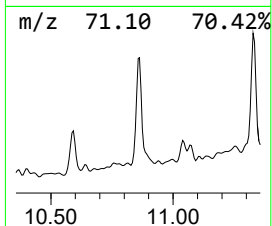
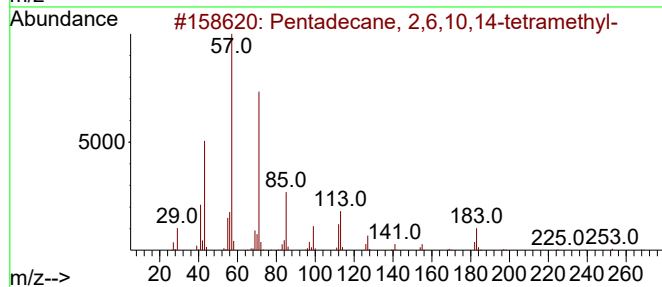
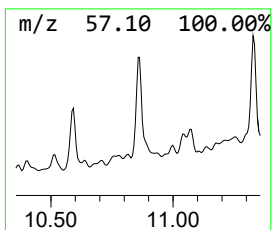
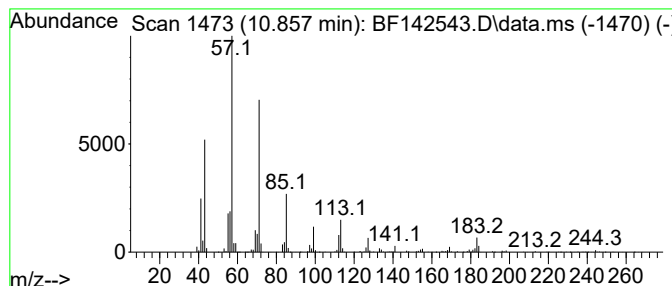
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TIC Integration Parameters: LSCINT.P

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Peak Number 18 Pentadecane, 2,6,10,14-tetr... Concentration Rank 2

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 10.857 | 20.00 ng | 1383130 | Phenanthrene-d10 | 11.433 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 95   |
| 2    |    |   | Pentadecane, 2,6,10-trimethyl-      | 254 | C18H38  | 003892-00-0 | 90   |
| 3    |    |   | Tridecane, 6-methyl-                | 198 | C14H30  | 013287-21-3 | 81   |
| 4    |    |   | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32  | 074645-98-0 | 72   |
| 5    |    |   | Hexadecane, 2,6,10-trimethyl-       | 268 | C19H40  | 055000-52-7 | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

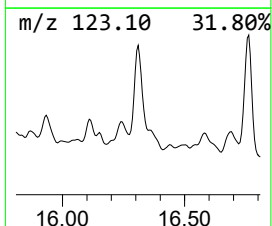
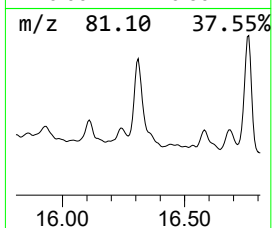
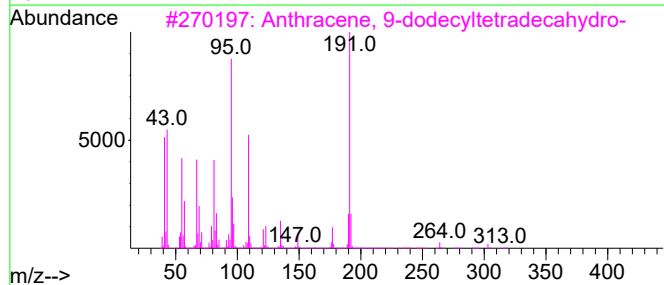
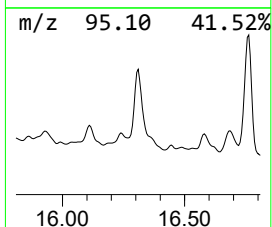
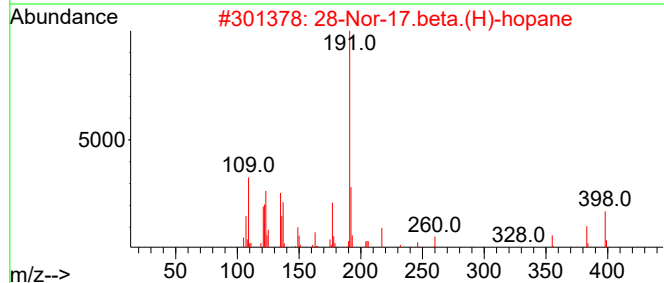
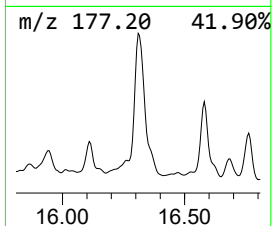
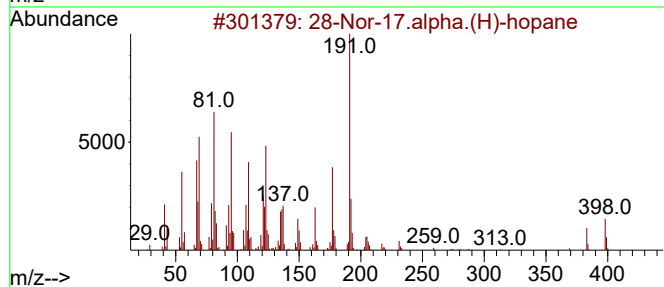
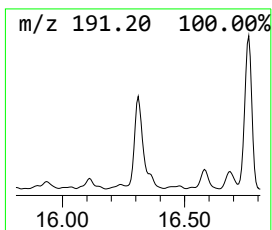
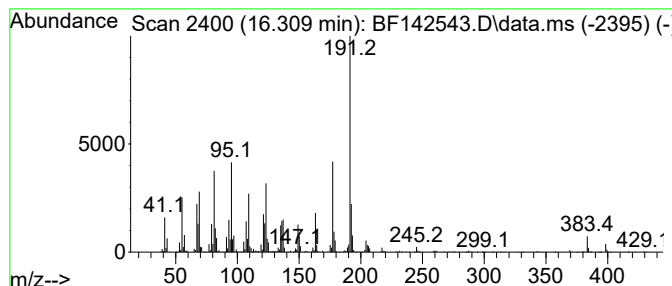
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 unknown16.310 Concentration Rank 3

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 16.310 | 20.00 ng | 1885100 | Perylene-d12     | 15.586 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|------|-------------------------------------|-----|----------|--------------|------|
| 1    |      | 28-Nor-17.alpha.(H)-hopane          | 398 | C29H50   | 053584-60-4  | 46   |
| 2    |      | 28-Nor-17.beta.(H)-hopane           | 398 | C29H50   | 036728-72-0  | 38   |
| 3    |      | Anthracene, 9-dodecyltetradecahy... | 360 | C26H48   | 055401-75-7  | 38   |
| 4    |      | .beta.-iso-Methyl ionone            | 206 | C14H22O  | 1000285-40-2 | 38   |
| 5    |      | (7S,7aR,13aS)-9-Methoxy-4,4,7,7a... | 344 | C22H32O3 | 1000482-45-3 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

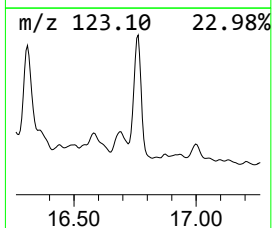
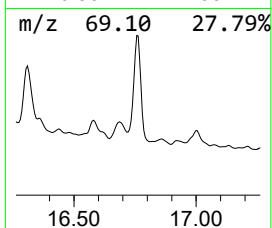
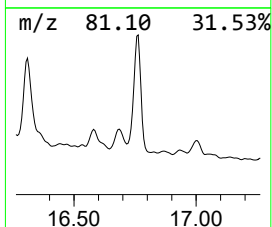
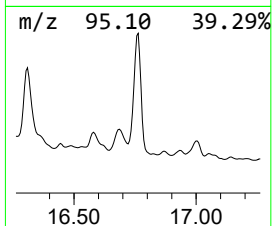
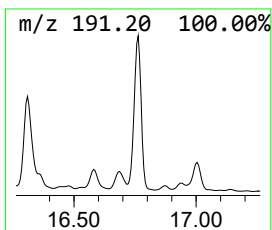
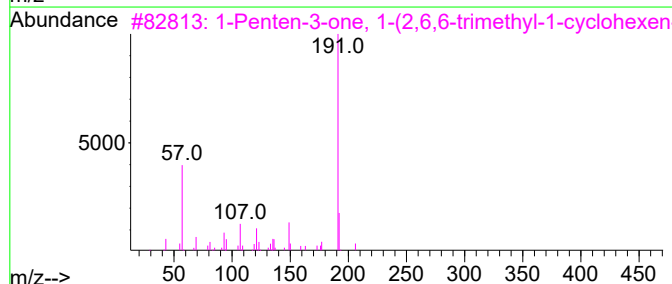
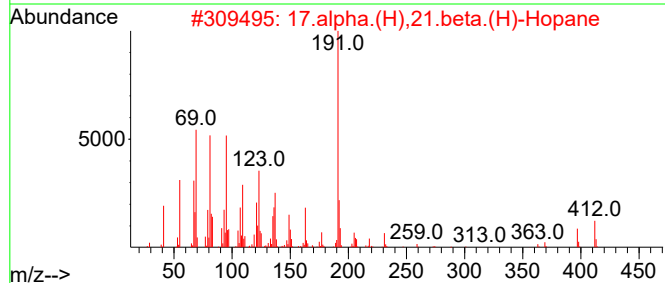
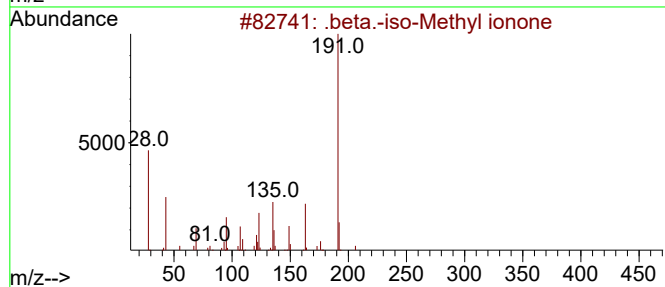
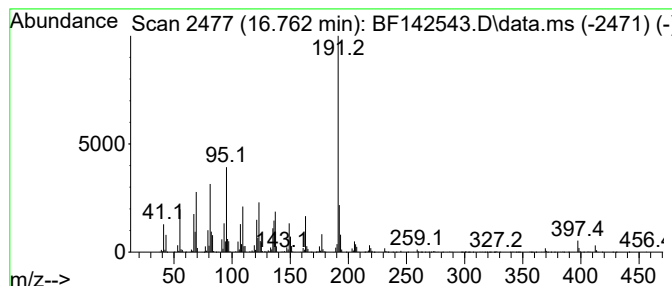
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 20 .beta.-iso-Methyl ionone Concentration Rank 4

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 16.762 | 19.46 ng | 1834050 | Perylene-d12     | 15.586 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|------|-------------------------------------|-----|---------|--------------|------|
| 1    |      | .beta.-iso-Methyl ionone            | 206 | C14H22O | 1000285-40-2 | 64   |
| 2    |      | 17.alpha.(H),21.beta.(H)-Hopane     | 412 | C30H52  | 013849-96-2  | 60   |
| 3    |      | 1-Penten-3-one, 1-(2,6,6-trimeth... | 206 | C14H22O | 000127-43-5  | 60   |
| 4    |      | 28-Nor-17.alpha.(H)-hopane          | 398 | C29H50  | 053584-60-4  | 53   |
| 5    |      | 1-Cyclohexene, 1,3,3-trimethyl-2... | 206 | C14H22O | 1000197-08-4 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052325\  
Data File : BF142543.D  
Acq On : 23 May 2025 18:58  
Operator : RC/JU  
Sample : Q2102-03  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LAW-25-0078

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.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 |
| 2-Hexanone, 3-h... | 4.416  | 5.0     | ng    | 158720   | 1                     | 6.898  | 638294  | 20.0 |
| 4-Pentene-2-ol,... | 4.545  | 11.3    | ng    | 360098   | 1                     | 6.898  | 638294  | 20.0 |
| 2-Pentanone, 4-... | 5.104  | 5.8     | ng    | 183352   | 1                     | 6.898  | 638294  | 20.0 |
| Acetic acid, cy... | 5.381  | 12.4    | ng    | 394812   | 1                     | 6.898  | 638294  | 20.0 |
| 2,4-Dimethyl-2,... | 6.410  | 12.7    | ng    | 404733   | 1                     | 6.898  | 638294  | 20.0 |
| Butane, 1,3-dim... | 6.704  | 5.7     | ng    | 180187   | 1                     | 6.898  | 638294  | 20.0 |
| 1H-Pyrazole, 4,... | 6.975  | 160.4   | ng    | 5118000  | 1                     | 6.898  | 638294  | 20.0 |
| unknown7.045       | 7.045  | 3.2     | ng    | 102387   | 1                     | 6.898  | 638294  | 20.0 |
| unknown7.092       | 7.092  | 5.1     | ng    | 162164   | 1                     | 6.898  | 638294  | 20.0 |
| Cyclopropane, 1... | 7.134  | 3.0     | ng    | 96140    | 1                     | 6.898  | 638294  | 20.0 |
| unknown7.522       | 7.522  | 3.4     | ng    | 107344   | 1                     | 6.898  | 638294  | 20.0 |
| unknown7.651       | 7.651  | 9.8     | ng    | 623412   | 2                     | 8.181  | 1276620 | 20.0 |
| 2(3H)-Furanone,... | 7.745  | 3.9     | ng    | 249830   | 2                     | 8.181  | 1276620 | 20.0 |
| Cyclohexane, me... | 8.069  | 5.2     | ng    | 333253   | 2                     | 8.181  | 1276620 | 20.0 |
| unknown8.292       | 8.292  | 3.8     | ng    | 244463   | 2                     | 8.181  | 1276620 | 20.0 |
| unknown8.681       | 8.681  | 13.4    | ng    | 853637   | 2                     | 8.181  | 1276620 | 20.0 |
| Nonane, 5-butyl-   | 9.651  | 6.3     | ng    | 319061   | 3                     | 9.939  | 1011900 | 20.0 |
| Pentadecane, 2,... | 10.857 | 20.0    | ng    | 1383130  | 4                     | 11.433 | 1383130 | 20.0 |
| unknown16.310      | 16.310 | 20.0    | ng    | 1885100  | 6                     | 15.586 | 1885100 | 20.0 |
| .beta.-iso-Meth... | 16.762 | 19.5    | ng    | 1834050  | 6                     | 15.586 | 1885100 | 20.0 |