

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091225\  
 Data File : BF143731.D  
 Acq On : 12 Sep 2025 14:13  
 Operator : RC/JU  
 Sample : Q3075-08  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MOO-25-0260

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-BF091025.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF143731.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         | 5.104       | 488           | 495         | 501          | rBB2     | 25105          | 36964         | 1.69%           | 0.352%        |
| 2         | 5.504       | 557           | 563         | 568          | rBV      | 311518         | 395286        | 18.03%          | 3.761%        |
| 3         | 6.510       | 728           | 734         | 739          | rBV      | 285909         | 371474        | 16.94%          | 3.534%        |
| 4         | 6.893       | 794           | 799         | 804          | rBB      | 162579         | 196477        | 8.96%           | 1.869%        |
| 5         | 7.451       | 888           | 894         | 898          | rBV      | 197184         | 243574        | 11.11%          | 2.318%        |
| 6         | 8.175       | 1012          | 1017        | 1024         | rBV      | 211268         | 264162        | 12.05%          | 2.513%        |
| 7         | 8.998       | 1153          | 1157        | 1160         | rBV2     | 23466          | 39303         | 1.79%           | 0.374%        |
| 8         | 9.251       | 1195          | 1200        | 1204         | rBV      | 429218         | 551976        | 25.18%          | 5.252%        |
| 9         | 9.334       | 1210          | 1214        | 1218         | rBV3     | 77439          | 124153        | 5.66%           | 1.181%        |
| 10        | 9.651       | 1264          | 1268        | 1271         | rBV2     | 215421         | 358995        | 16.37%          | 3.416%        |
| 11        | 9.804       | 1291          | 1294        | 1297         | rBV4     | 171166         | 268264        | 12.24%          | 2.552%        |
| 12        | 9.851       | 1299          | 1302        | 1306         | rVB2     | 293976         | 380615        | 17.36%          | 3.621%        |
| 13        | 9.934       | 1311          | 1316        | 1321         | rVV4     | 340533         | 678394        | 30.94%          | 6.455%        |
| 14        | 10.357      | 1385          | 1388        | 1392         | rBV2     | 277313         | 346056        | 15.78%          | 3.293%        |
| 15        | 10.881      | 1474          | 1477        | 1483         | rVB      | 756255         | 992364        | 45.26%          | 9.442%        |
| 16        | 14.086      | 2019          | 2022        | 2025         | rBV3     | 309389         | 483782        | 22.07%          | 4.603%        |
| 17        | 15.139      | 2194          | 2201        | 2210         | rBV      | 864506         | 2192528       | 100.00%         | 20.861%       |
| 18        | 15.239      | 2215          | 2218        | 2223         | rVB      | 109045         | 141828        | 6.47%           | 1.349%        |
| 19        | 15.445      | 2247          | 2253        | 2258         | rVB      | 460323         | 784314        | 35.77%          | 7.462%        |
| 20        | 15.510      | 2258          | 2264        | 2269         | rBV      | 388956         | 622626        | 28.40%          | 5.924%        |
| 21        | 15.563      | 2269          | 2273        | 2277         | rBV2     | 168613         | 291245        | 13.28%          | 2.771%        |
| 22        | 17.057      | 2521          | 2527        | 2535         | rVB2     | 180641         | 388902        | 17.74%          | 3.700%        |
| 23        | 17.239      | 2553          | 2558        | 2562         | rBV      | 26120          | 59058         | 2.69%           | 0.562%        |
| 24        | 17.510      | 2597          | 2604        | 2617         | rVB      | 120336         | 297801        | 13.58%          | 2.833%        |

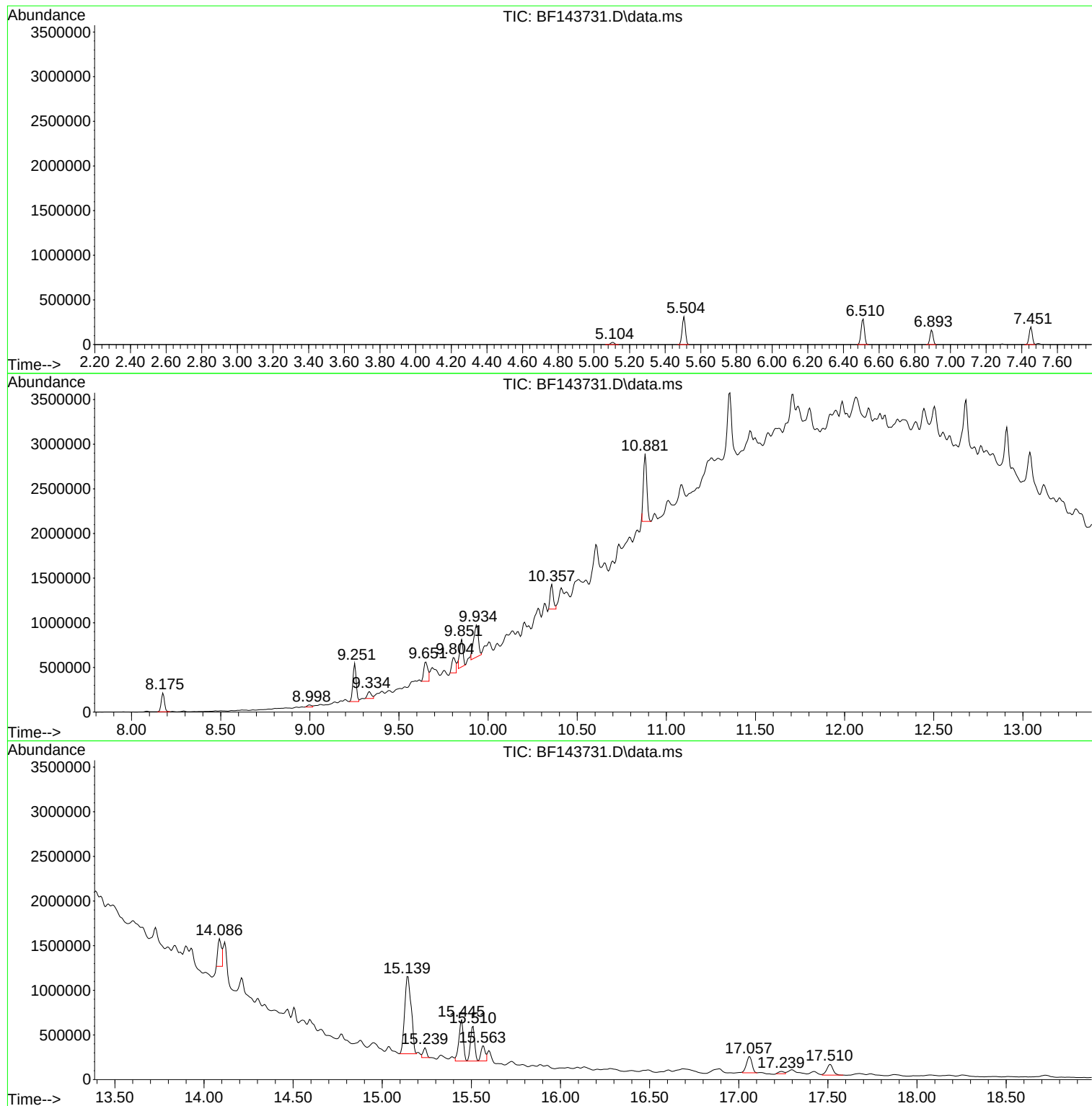
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091225\  
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Sample : Q3075-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-25-0260

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091025.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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BNA\_F  
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MOO-25-0260

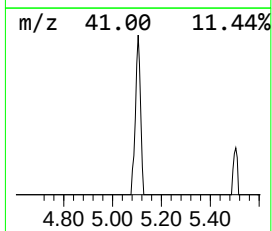
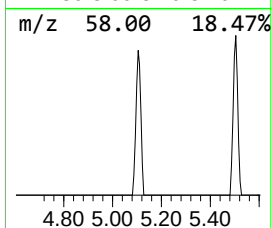
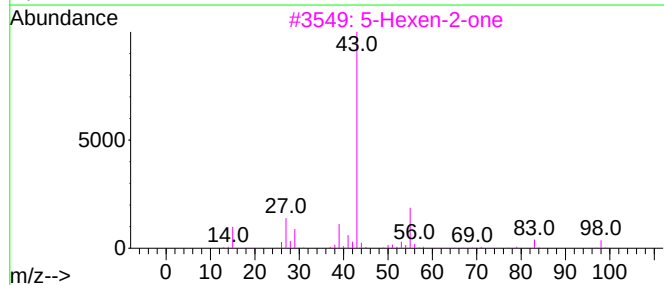
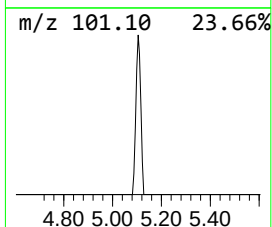
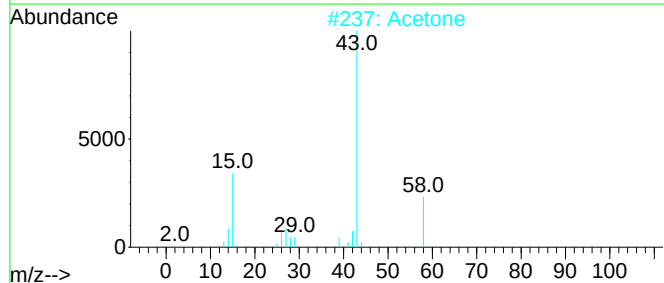
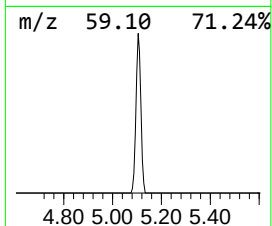
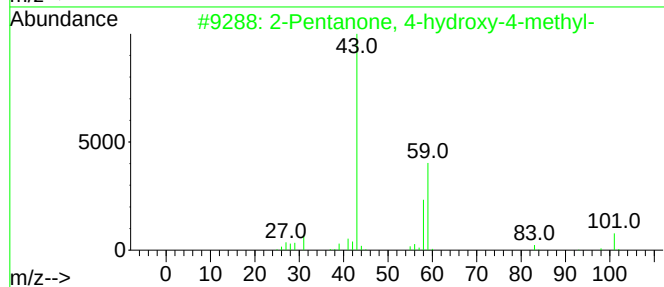
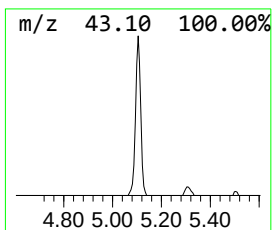
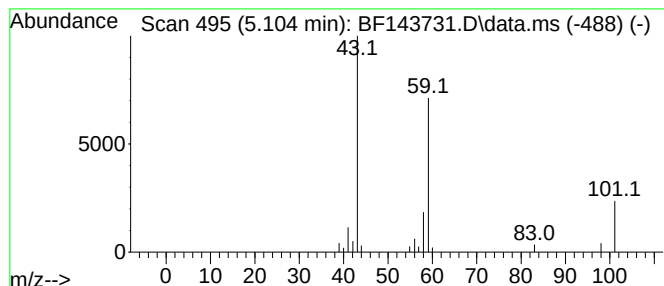
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD       | R.T.  |
|-------|---------|-------|------------------------|-------|
| 5.104 | 3.76 ng | 36964 | 1,4-Dichlorobenzene-d4 | 6.893 |

| Hit# | of 5                             | Tentative ID | MW      | MolForm     | CAS# | Qual |
|------|----------------------------------|--------------|---------|-------------|------|------|
| 1    | 2-Pentanone, 4-hydroxy-4-methyl- | 116          | C6H12O2 | 000123-42-2 | 56   |      |
| 2    | Acetone                          | 58           | C3H6O   | 000067-64-1 | 9    |      |
| 3    | 5-Hexen-2-one                    | 98           | C6H10O  | 000109-49-9 | 9    |      |
| 4    | 3-Hydroxy-3-methyl-2-butanone    | 102          | C5H10O2 | 000115-22-0 | 9    |      |
| 5    | 3-Octanol                        | 130          | C8H18O  | 000589-98-0 | 9    |      |



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Sample : Q3075-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

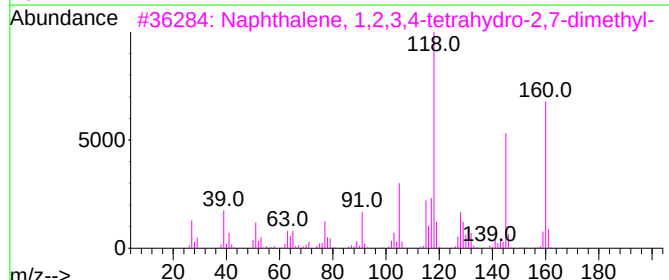
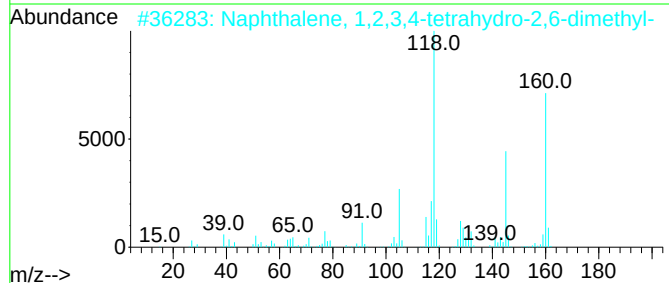
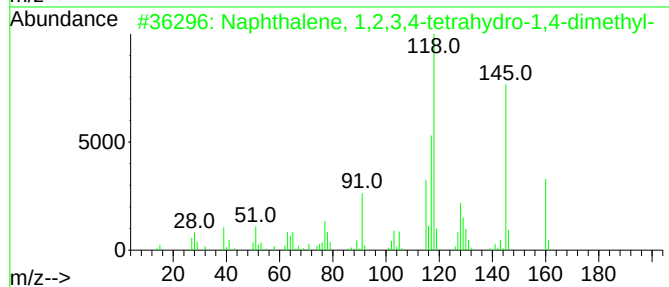
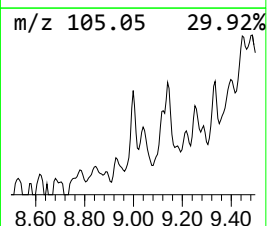
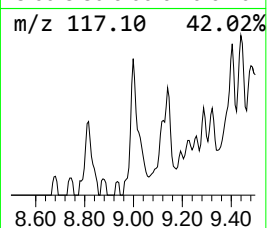
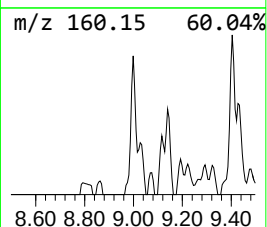
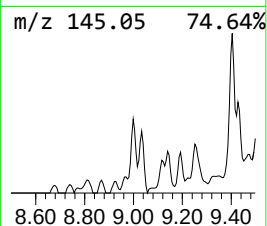
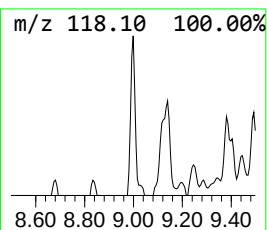
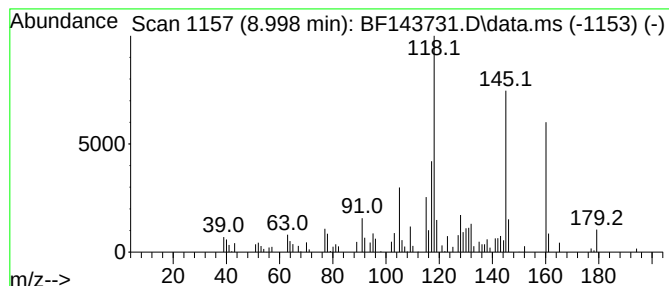
Instrument :  
BNA\_F  
ClientSampleId :  
MOO-25-0260

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091025.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 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 10

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|-------|------------------|-------------|------|
| 8.998     | 2.98 ng                             | 39303 | Naphthalene-d8   | 8.175       |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 1,2,3,4-tetrahydro-... | 160   | C12H16           | 004175-54-6 | 97   |
| 2         | Naphthalene, 1,2,3,4-tetrahydro-... | 160   | C12H16           | 007524-63-2 | 81   |
| 3         | Naphthalene, 1,2,3,4-tetrahydro-... | 160   | C12H16           | 013065-07-1 | 76   |
| 4         | 2,5-Dimethylbenzyl cyanide          | 145   | C10H11N          | 016213-85-7 | 58   |
| 5         | Benzene, 1-(1-methylethenyl)-3-(... | 160   | C12H16           | 001129-29-9 | 55   |



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Sample : Q3075-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-25-0260

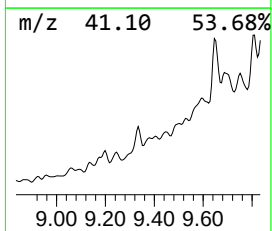
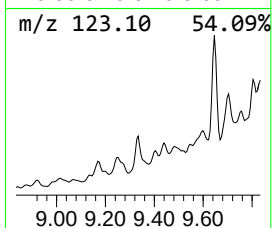
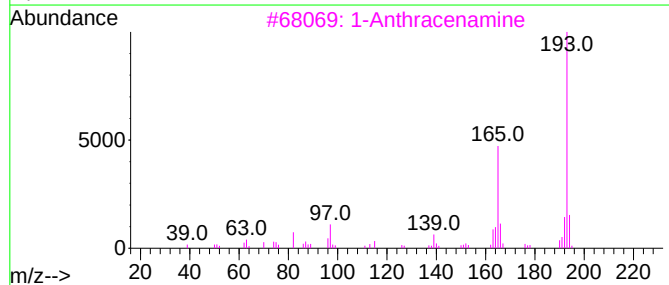
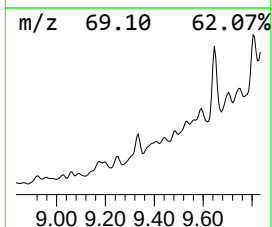
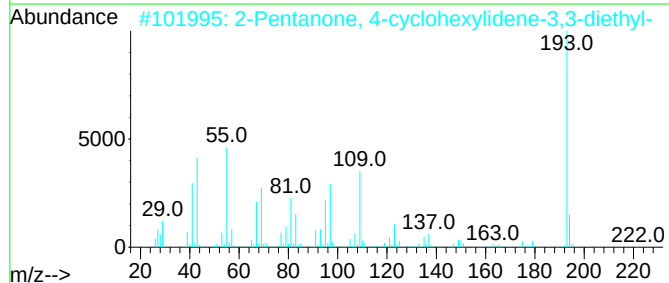
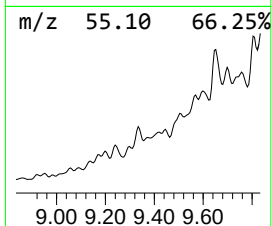
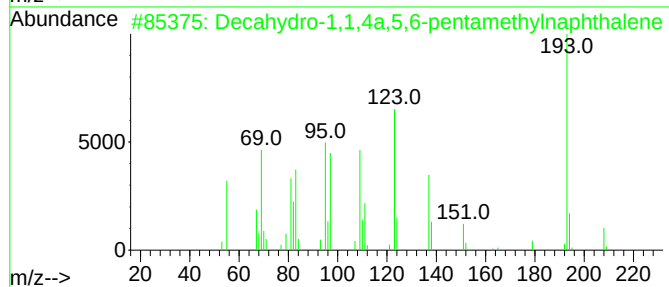
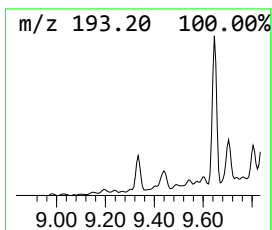
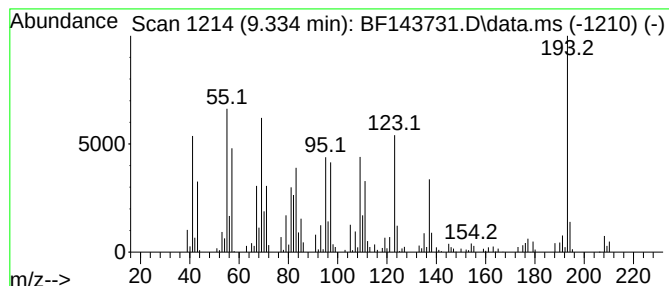
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.334 | 3.66 ng | 124153 | Acenaphthene-d10 | 9.939 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Decahydro-1,1,4a,5,6-pentamethyl... | 208 | C15H28  | 080655-44-3 | 96   |
| 2    |    |   | 2-Pentanone, 4-cyclohexylidene-3... | 222 | C15H26O | 313253-65-5 | 43   |
| 3    |    |   | 1-Anthracenamine                    | 193 | C14H11N | 000610-49-1 | 25   |
| 4    |    |   | 2-Anthracenamine                    | 193 | C14H11N | 000613-13-8 | 22   |
| 5    |    |   | 3-Phenylindole                      | 193 | C14H11N | 001504-16-1 | 18   |



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Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091025.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

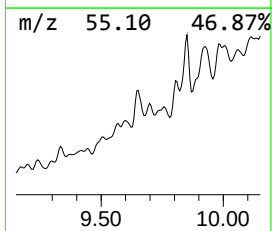
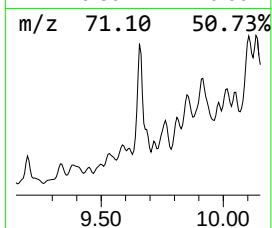
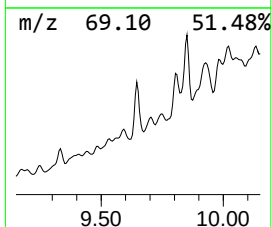
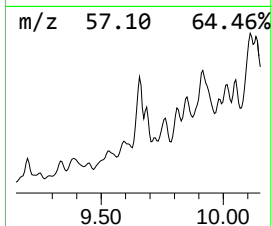
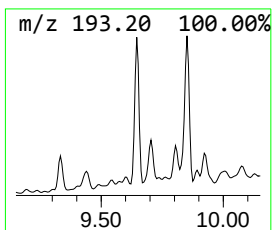
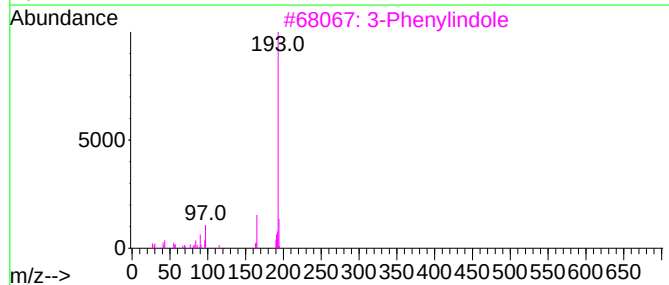
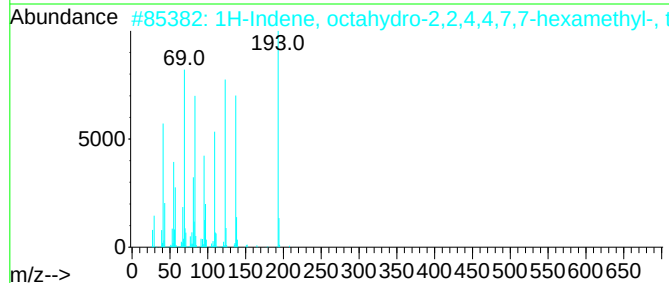
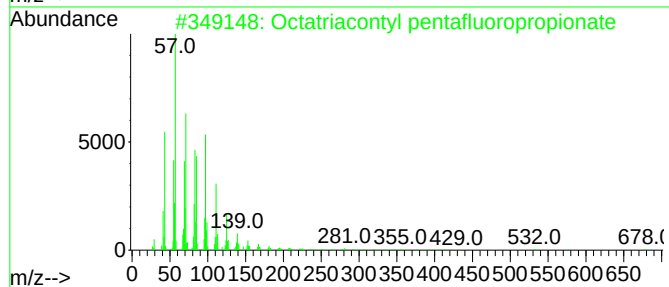
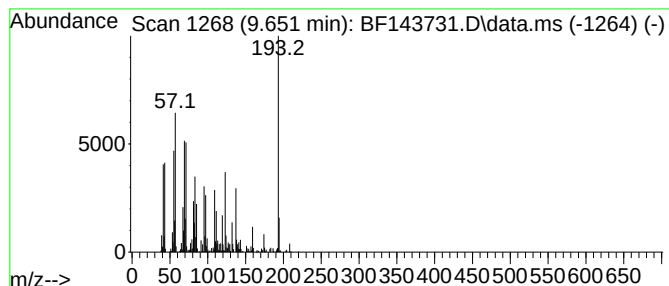
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown9.651 Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 9.651 | 10.58 ng | 358995 | Acenaphthene-d10 | 9.939 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Octatriacontyl pentafluoropropio... | 697 | C41H77F5O2 | 1000351-89-1 | 35   |
| 2    |    |   | 1H-Indene, octahydro-2,2,4,4,7,7... | 208 | C15H28     | 054832-83-6  | 35   |
| 3    |    |   | 3-Phenylindole                      | 193 | C14H11N    | 001504-16-1  | 27   |
| 4    |    |   | 2-((Trimethylsilyl)thio)nicotino... | 208 | C9H12N2SSi | 1000474-02-5 | 22   |
| 5    |    |   | Iminostilbene                       | 193 | C14H11N    | 000256-96-2  | 22   |



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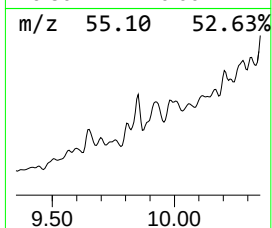
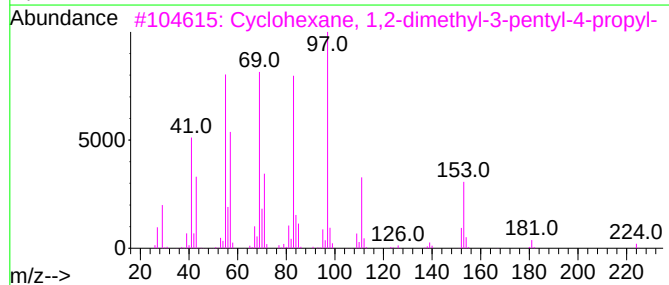
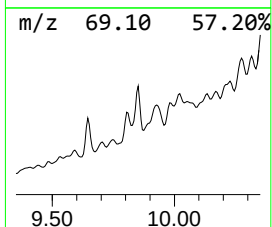
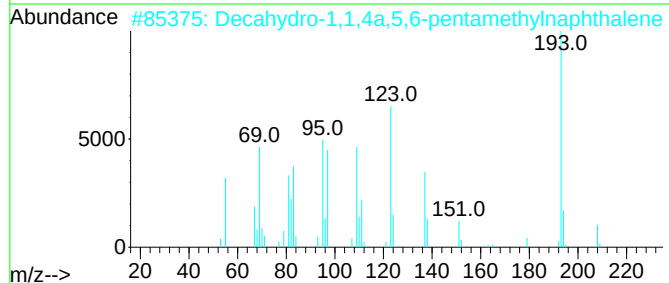
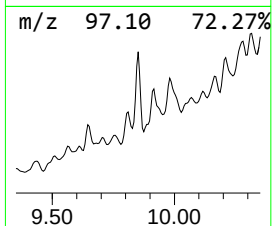
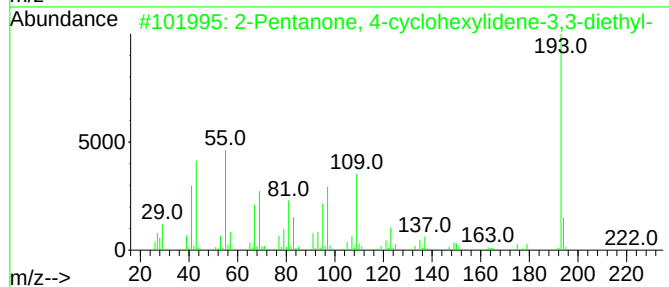
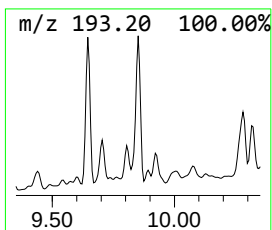
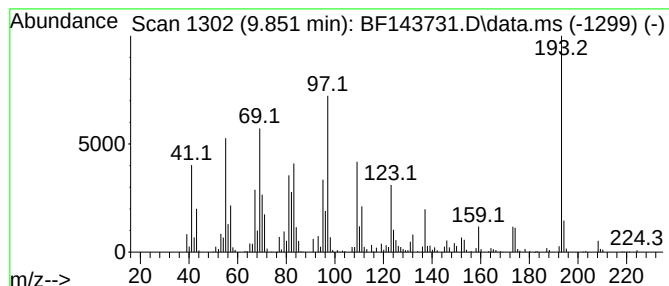
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 2-Pentanone, 4-cyclohexylid... Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 9.851 | 11.22 ng | 380615 | Acenaphthene-d10 | 9.939 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2-Pentanone, 4-cyclohexylidene-3... | 222 | C15H26O      | 313253-65-5  | 50      |      |      |
| 2    | Decahydro-1,1,4a,5,6-pentamethyl... | 208 | C15H28       | 080655-44-3  | 49      |      |      |
| 3    | Cyclohexane, 1,2-dimethyl-3-pent... | 224 | C16H32       | 062376-17-4  | 38      |      |      |
| 4    | Acetamide, N-methyl-N-[4-[3-(hyd... | 224 | C12H20N2O2   | 130403-63-3  | 18      |      |      |
| 5    | 2-((Trimethylsilyl)thio)nicotino... | 208 | C9H12N2SSi   | 1000474-02-5 | 18      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091225\  
Data File : BF143731.D  
Acq On : 12 Sep 2025 14:13  
Operator : RC/JU  
Sample : Q3075-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-25-0260

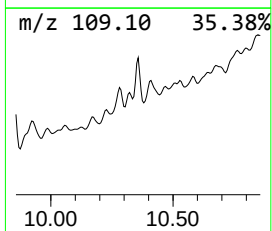
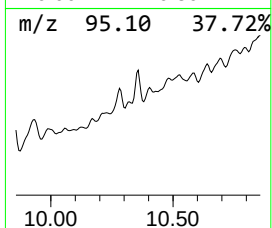
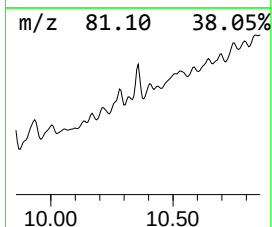
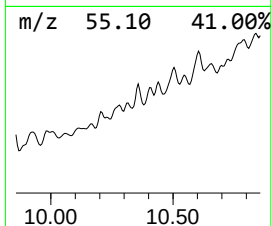
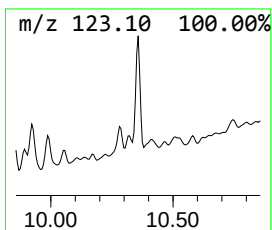
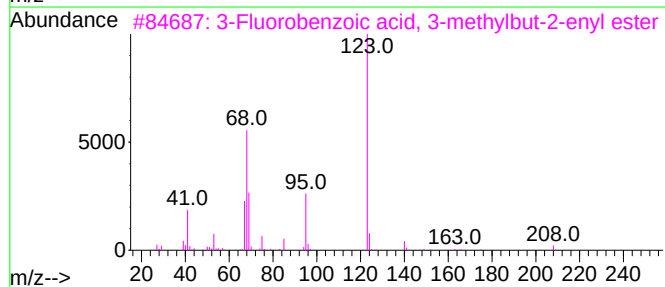
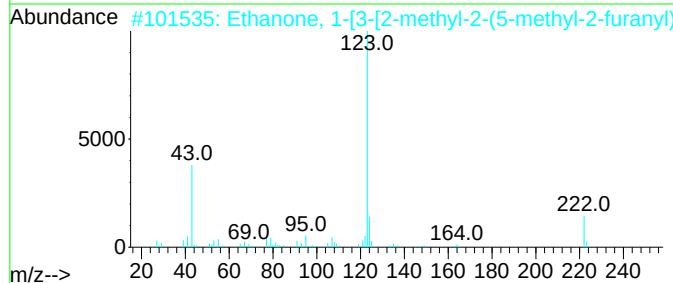
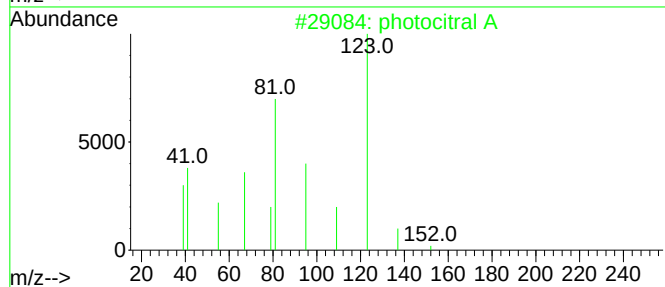
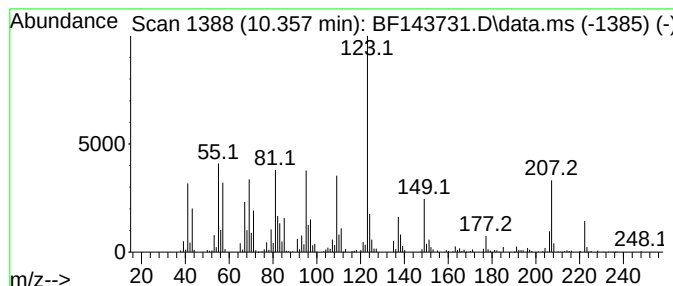
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 unknown10.357 Concentration Rank 5

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.  |
|--------|----------|--------|------------------|-------|
| 10.357 | 10.20 ng | 346056 | Acenaphthene-d10 | 9.939 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | photocitral A                       | 152 | C10H16O    | 055253-28-6  | 46   |
| 2    |    |   | Ethanone, 1-[3-[2-methyl-2-(5-me... | 222 | C13H18O3   | 080114-28-9  | 45   |
| 3    |    |   | 3-Fluorobenzoic acid, 3-methylbu... | 208 | C12H13FO2  | 154559-38-3  | 43   |
| 4    |    |   | (Phenylthio)acetic acid, cyclobu... | 222 | C12H14O2S  | 1010299-95-5 | 43   |
| 5    |    |   | Bicyclo[4.1.0]heptane-7-carboxam... | 222 | C11H14N2OS | 329912-72-3  | 43   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091225\  
Data File : BF143731.D  
Acq On : 12 Sep 2025 14:13  
Operator : RC/JU  
Sample : Q3075-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-25-0260

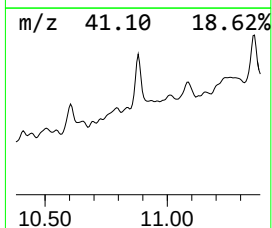
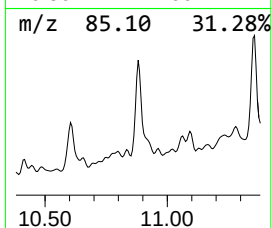
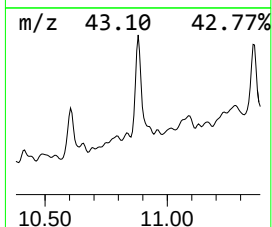
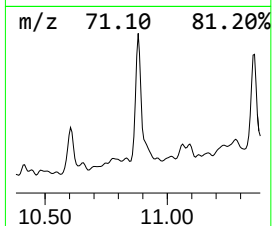
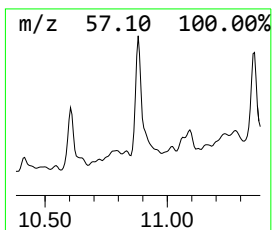
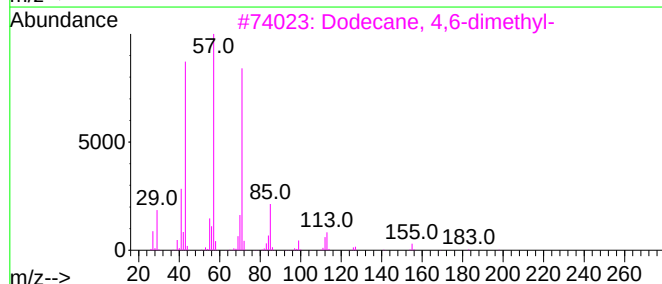
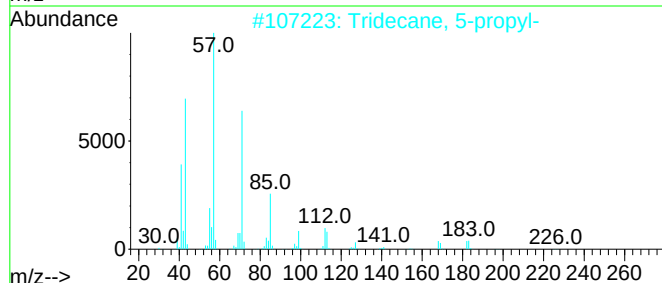
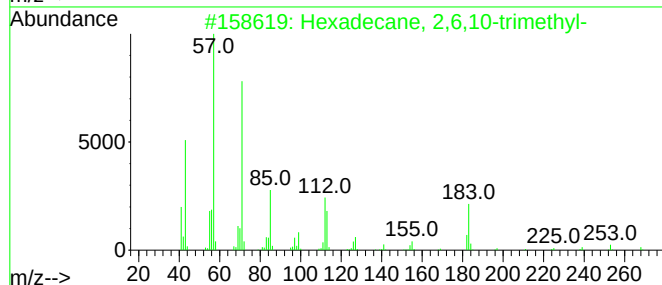
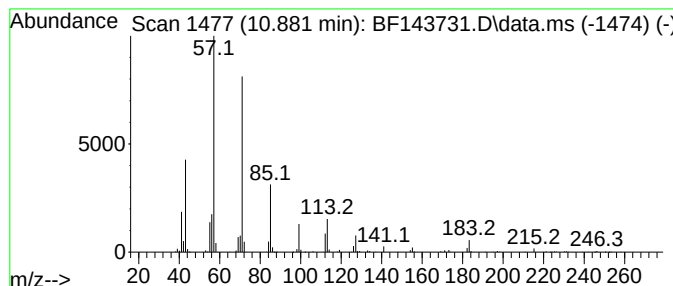
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Hexadecane, 2,6,10-trimethyl- Concentration Rank 2

| R.T.      | EstConc                        | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|--------------------------------|--------|------------------|---------|-------------|------|
| 10.881    | 20.00 ng                       | 992364 | Phenanthrene-d10 |         | 11.445      |      |
| Hit# of 5 | Tentative ID                   |        | MW               | MolForm | CAS#        | Qual |
| 1         | Hexadecane, 2,6,10-trimethyl-  |        | 268              | C19H40  | 055000-52-7 | 83   |
| 2         | Tridecane, 5-propyl-           |        | 226              | C16H34  | 055045-11-9 | 72   |
| 3         | Dodecane, 4,6-dimethyl-        |        | 198              | C14H30  | 061141-72-8 | 72   |
| 4         | Pentadecane, 2,6,10-trimethyl- |        | 254              | C18H38  | 003892-00-0 | 62   |
| 5         | 3-Bromooctane                  |        | 192              | C8H17Br | 000999-64-4 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091225\  
Data File : BF143731.D  
Acq On : 12 Sep 2025 14:13  
Operator : RC/JU  
Sample : Q3075-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-25-0260

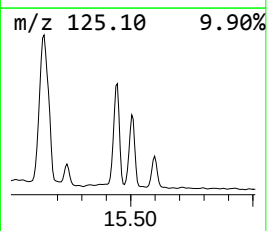
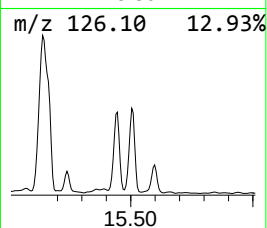
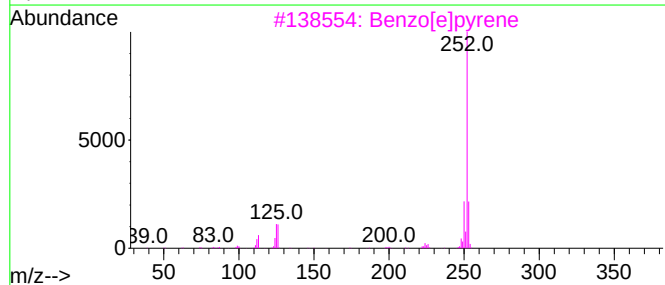
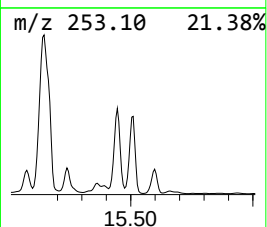
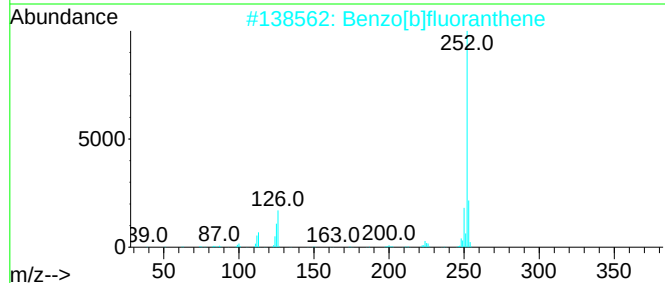
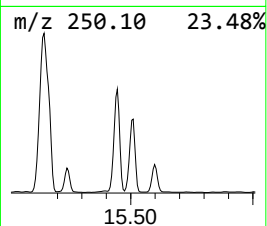
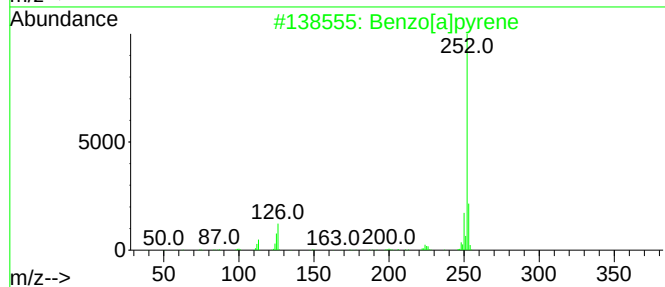
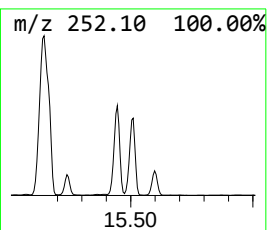
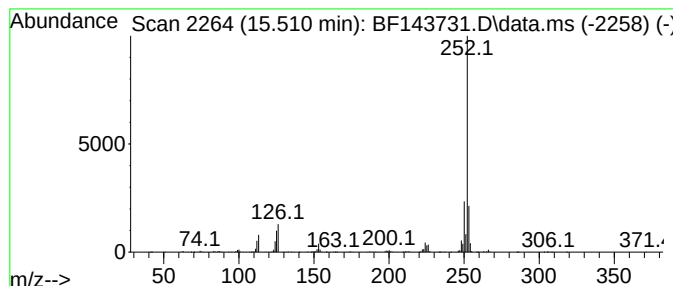
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Benzo[a]pyrene Concentration Rank 1

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 15.510 | 42.76 ng | 622626 | Perylene-d12     | 15.569 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[a]pyrene       | 252 | C20H12  | 000050-32-8 | 98   |
| 2    |    |   | Benzo[b]fluoranthene | 252 | C20H12  | 000205-99-2 | 97   |
| 3    |    |   | Benzo[e]pyrene       | 252 | C20H12  | 000192-97-2 | 96   |
| 4    |    |   | Perylene             | 252 | C20H12  | 000198-55-0 | 94   |
| 5    |    |   | Benzo[k]fluoranthene | 252 | C20H12  | 000207-08-9 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091225\  
Data File : BF143731.D  
Acq On : 12 Sep 2025 14:13  
Operator : RC/JU  
Sample : Q3075-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-25-0260

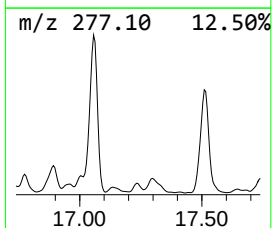
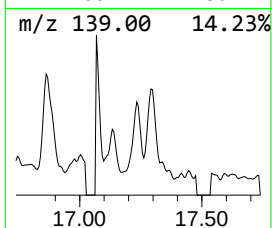
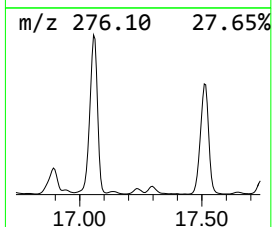
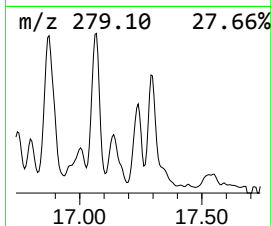
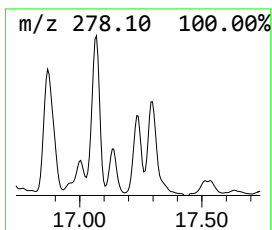
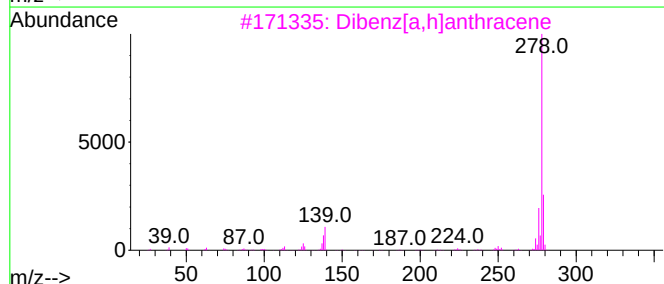
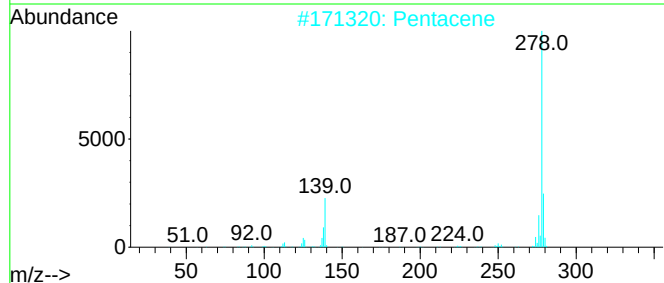
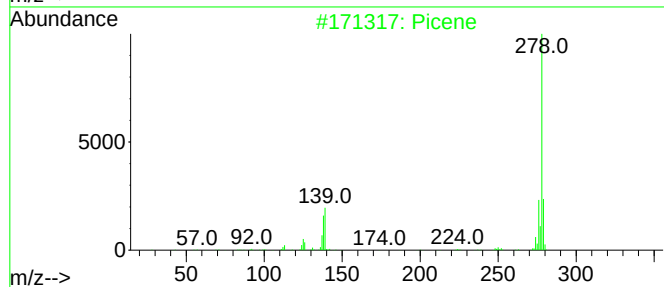
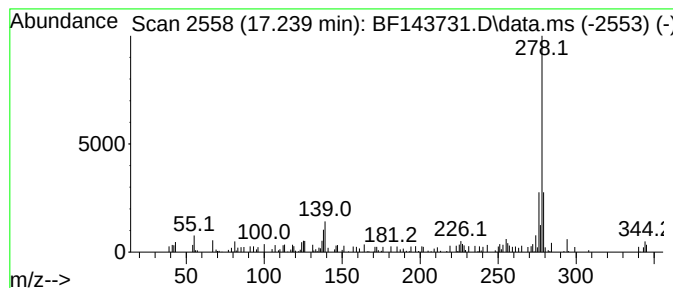
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Picene Concentration Rank 7

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 17.239 | 4.06 ng | 59058 | Perylene-d12     | 15.569 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Picene                              | 278 | C22H14  | 000213-46-7 | 91   |
| 2    |      | Pentacene                           | 278 | C22H14  | 000135-48-8 | 91   |
| 3    |      | Dibenz[a,h]anthracene               | 278 | C22H14  | 000053-70-3 | 83   |
| 4    |      | Vanadium, (.eta.5-2,4-cyclopenta... | 278 | C17H23V | 126948-97-8 | 72   |
| 5    |      | Benzo[b]triphenylene                | 278 | C22H14  | 000215-58-7 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091225\  
Data File : BF143731.D  
Acq On : 12 Sep 2025 14:13  
Operator : RC/JU  
Sample : Q3075-08  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-25-0260

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091025.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-Pentanone, 4-... | 5.104  | 3.8     | ng    | 36964    | 1                     | 6.893  | 196477 | 20.0 |
| Naphthalene, 1,... | 8.998  | 3.0     | ng    | 39303    | 2                     | 8.175  | 264162 | 20.0 |
| Decahydro-1,1,4... | 9.334  | 3.7     | ng    | 124153   | 3                     | 9.939  | 678394 | 20.0 |
| unknown9.651       | 9.651  | 10.6    | ng    | 358995   | 3                     | 9.939  | 678394 | 20.0 |
| 2-Pentanone, 4-... | 9.851  | 11.2    | ng    | 380615   | 3                     | 9.939  | 678394 | 20.0 |
| unknown10.357      | 10.357 | 10.2    | ng    | 346056   | 3                     | 9.939  | 678394 | 20.0 |
| Hexadecane, 2,6... | 10.881 | 20.0    | ng    | 992364   | 4                     | 11.445 | 992364 | 20.0 |
| Benzo[a]pyrene     | 15.510 | 42.8    | ng    | 622626   | 6                     | 15.569 | 291245 | 20.0 |
| Picene             | 17.239 | 4.1     | ng    | 59058    | 6                     | 15.569 | 291245 | 20.0 |