

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
 Data File : BF141040.D  
 Acq On : 30 Dec 2024 19:25  
 Operator : RC/JU  
 Sample : P5386-03  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 MOO-24-00395-96

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-BF122624.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF141040.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 2.399       | 46            | 52          | 60           | rVB      | 217826         | 357977        | 6.01%           | 0.497%        |
| 2         | 5.434       | 558           | 568         | 577          | rBV      | 3545882        | 4858920       | 81.56%          | 6.740%        |
| 3         | 6.446       | 733           | 740         | 746          | rBV      | 3413892        | 4856777       | 81.53%          | 6.737%        |
| 4         | 6.828       | 800           | 805         | 810          | rBV      | 1123455        | 1400971       | 23.52%          | 1.943%        |
| 5         | 6.945       | 820           | 825         | 831          | rBV      | 217519         | 289766        | 4.86%           | 0.402%        |
| 6         | 7.387       | 892           | 900         | 905          | rBV      | 2277248        | 3012278       | 50.56%          | 4.179%        |
| 7         | 7.892       | 981           | 986         | 990          | rBV      | 654960         | 815554        | 13.69%          | 1.131%        |
| 8         | 8.104       | 1016          | 1022        | 1030         | rBV      | 1394535        | 1845793       | 30.98%          | 2.561%        |
| 9         | 8.704       | 1119          | 1124        | 1126         | rBV3     | 146867         | 215543        | 3.62%           | 0.299%        |
| 10        | 8.822       | 1137          | 1144        | 1148         | rBV2     | 193400         | 316558        | 5.31%           | 0.439%        |
| 11        | 9.063       | 1181          | 1185        | 1188         | rBV3     | 156854         | 239427        | 4.02%           | 0.332%        |
| 12        | 9.181       | 1200          | 1205        | 1209         | rBV      | 4064987        | 5150582       | 86.46%          | 7.145%        |
| 13        | 9.257       | 1214          | 1218        | 1223         | rBV2     | 961500         | 1698929       | 28.52%          | 2.357%        |
| 14        | 9.498       | 1255          | 1259        | 1261         | rBV3     | 467798         | 758186        | 12.73%          | 1.052%        |
| 15        | 9.575       | 1269          | 1272        | 1277         | rBV3     | 652062         | 1420970       | 23.85%          | 1.971%        |
| 16        | 9.681       | 1287          | 1290        | 1294         | rVB4     | 598874         | 710217        | 11.92%          | 0.985%        |
| 17        | 9.739       | 1296          | 1300        | 1302         | rBV3     | 894421         | 1521565       | 25.54%          | 2.111%        |
| 18        | 9.857       | 1315          | 1320        | 1324         | rBV2     | 1749393        | 3370998       | 56.59%          | 4.676%        |
| 19        | 13.998      | 2020          | 2024        | 2031         | rVB      | 1805618        | 2864002       | 48.08%          | 3.973%        |
| 20        | 14.127      | 2039          | 2046        | 2055         | rVB2     | 2721702        | 5744367       | 96.42%          | 7.969%        |
| 21        | 14.639      | 2128          | 2133        | 2142         | rVB3     | 812785         | 1926330       | 32.34%          | 2.672%        |
| 22        | 14.768      | 2149          | 2155        | 2164         | rVB2     | 3049755        | 5836671       | 97.97%          | 8.097%        |
| 23        | 15.363      | 2252          | 2256        | 2265         | rVB2     | 565357         | 972643        | 16.33%          | 1.349%        |
| 24        | 15.445      | 2266          | 2270        | 2274         | rVV      | 959856         | 1399531       | 23.49%          | 1.941%        |
| 25        | 15.521      | 2277          | 2283        | 2294         | rVB      | 2671536        | 5957348       | 100.00%         | 8.264%        |
| 26        | 16.310      | 2412          | 2417        | 2423         | rBV      | 376359         | 684622        | 11.49%          | 0.950%        |
| 27        | 16.410      | 2429          | 2434        | 2440         | rBV      | 457440         | 855807        | 14.37%          | 1.187%        |
| 28        | 16.521      | 2446          | 2453        | 2459         | rBV      | 2097134        | 4746830       | 79.68%          | 6.585%        |
| 29        | 17.498      | 2613          | 2619        | 2629         | rBV2     | 824123         | 2029581       | 34.07%          | 2.815%        |
| 30        | 17.604      | 2630          | 2637        | 2655         | rVB      | 762326         | 2228590       | 37.41%          | 3.092%        |
| 31        | 17.909      | 2679          | 2689        | 2697         | rBV2     | 1324882        | 3999804       | 67.14%          | 5.549%        |

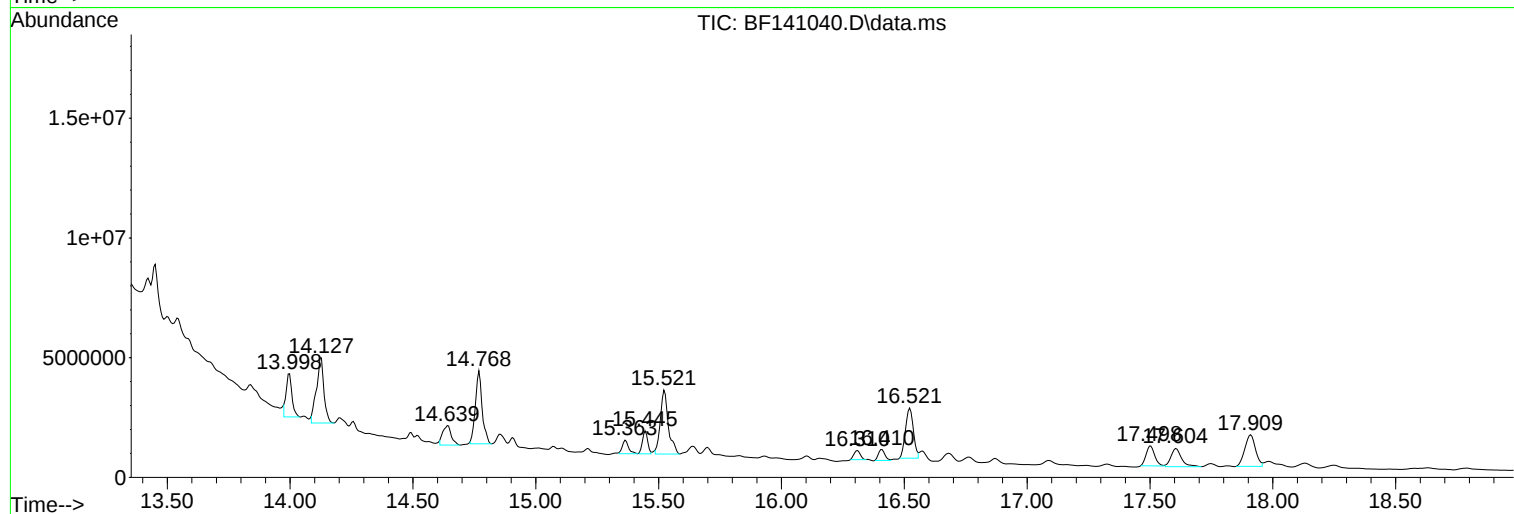
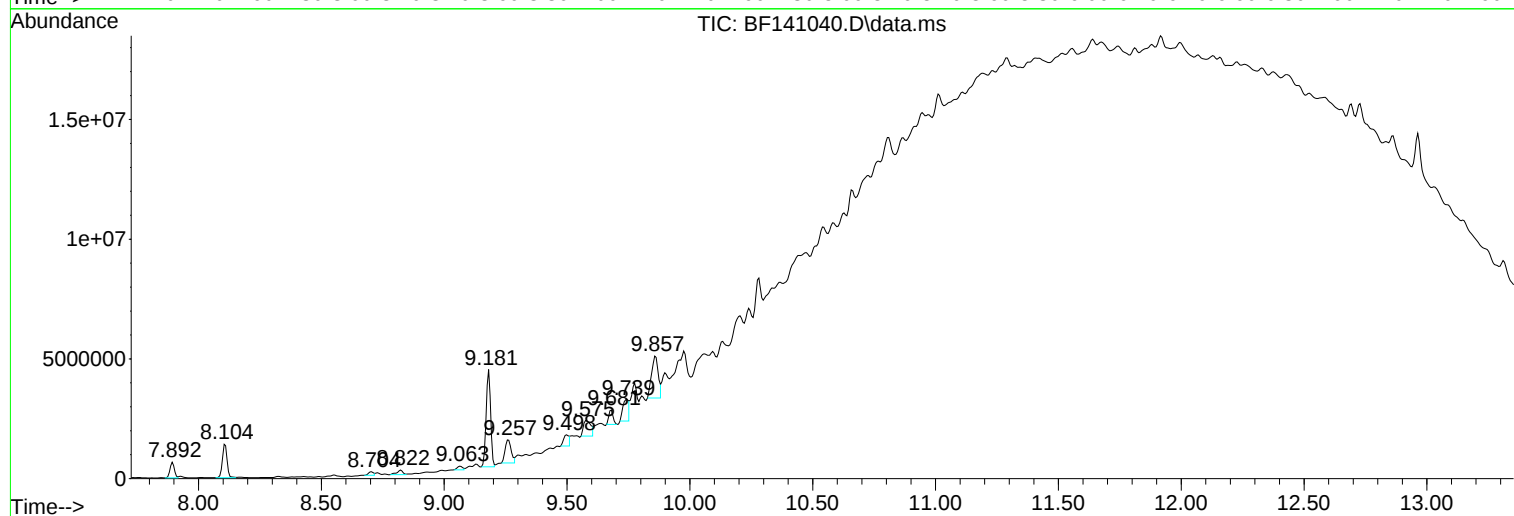
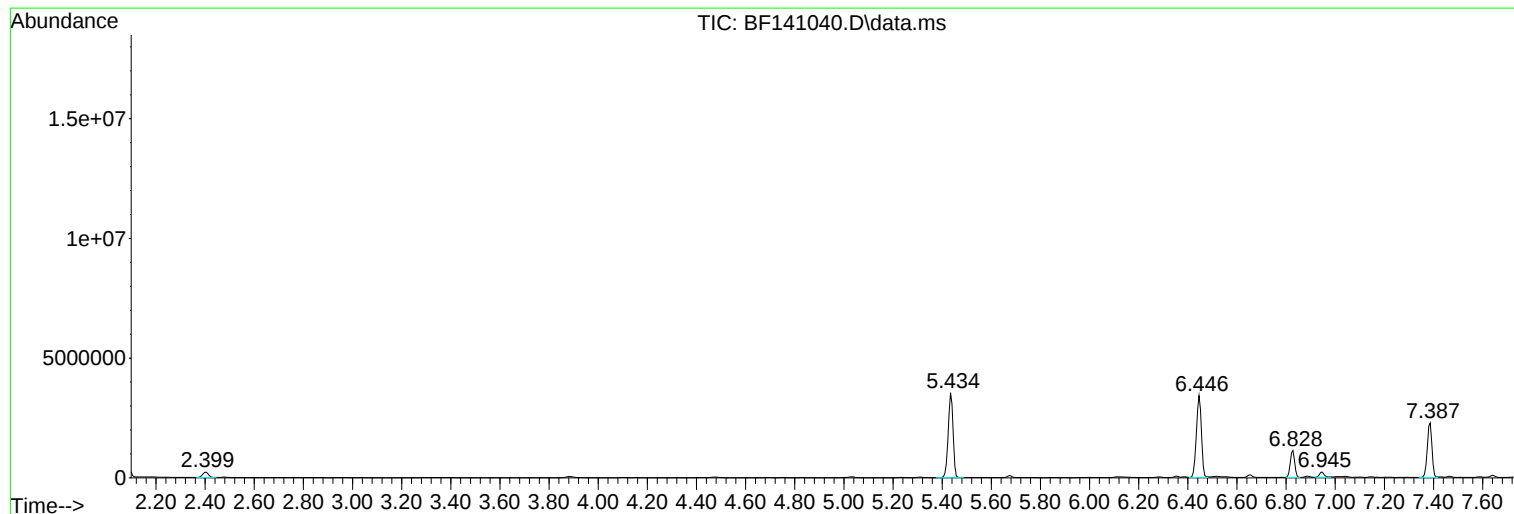
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Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.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\BF123024\  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

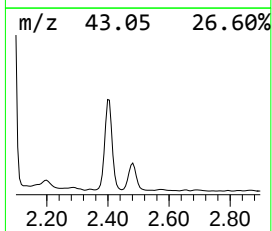
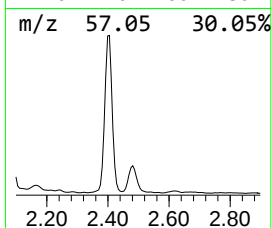
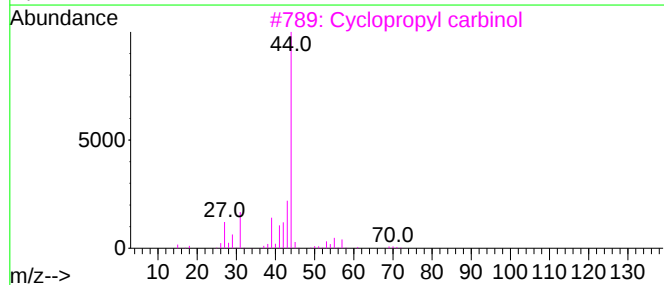
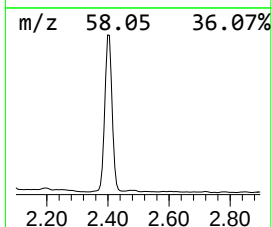
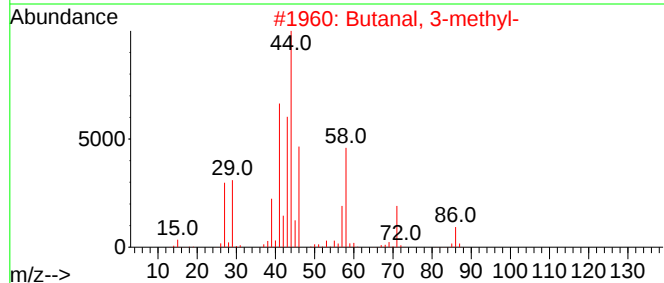
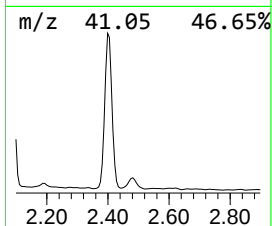
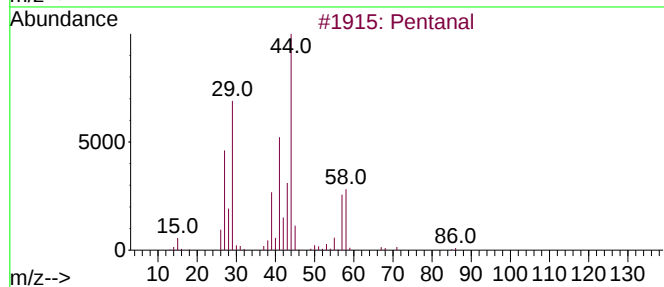
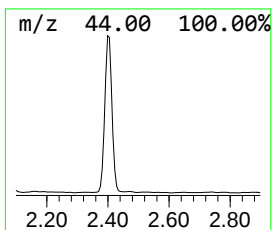
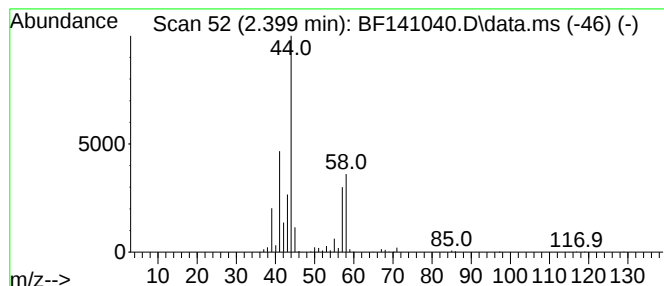
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Pentanal Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 2.399 | 5.11 ng | 357977 | 1,4-Dichlorobenzene-d4 | 6.828 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentanal                 | 86  | C5H10O  | 000110-62-3 | 83   |
| 2    |    |   | Butanal, 3-methyl-       | 86  | C5H10O  | 000590-86-3 | 64   |
| 3    |    |   | Cyclopropyl carbinol     | 72  | C4H8O   | 002516-33-8 | 36   |
| 4    |    |   | Glycidol                 | 74  | C3H6O2  | 000556-52-5 | 36   |
| 5    |    |   | 2-Heptanamine, 5-methyl- | 129 | C8H19N  | 053907-81-6 | 9    |



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Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

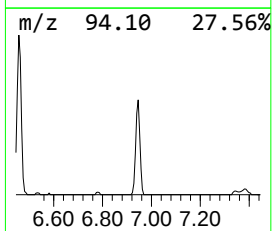
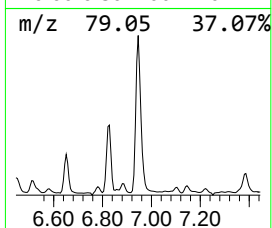
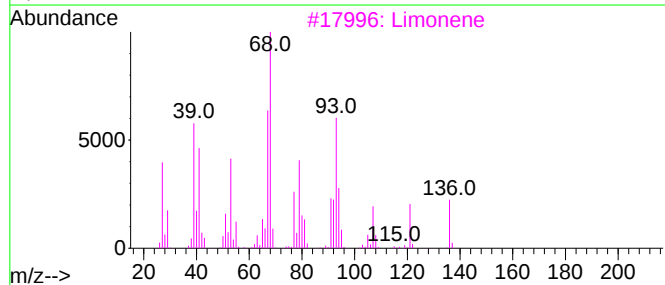
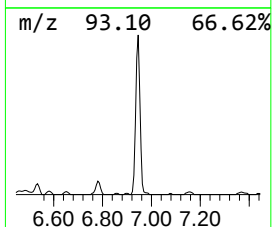
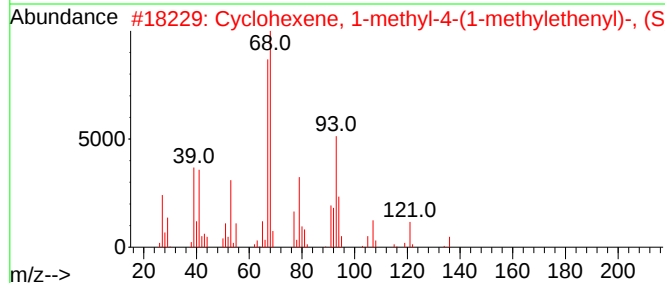
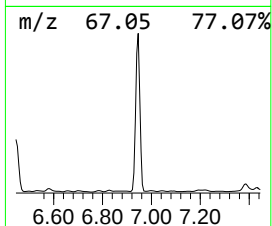
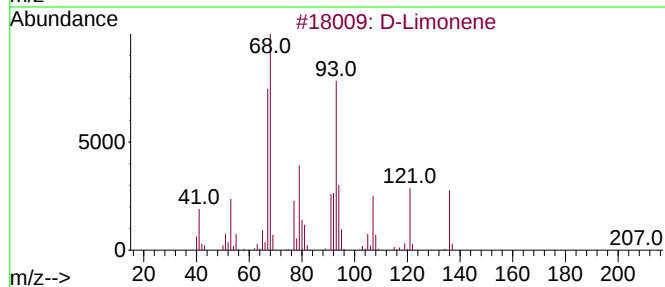
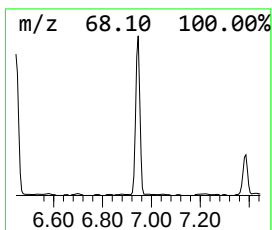
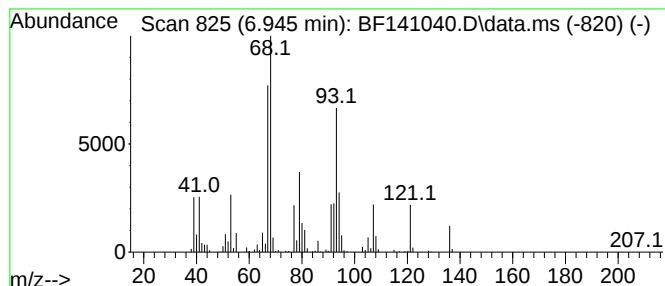
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 D-Limonene Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.945 | 4.14 ng | 289766 | 1,4-Dichlorobenzene-d4 | 6.828 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | D-Limonene                          | 136 | C10H16   | 005989-27-5 | 98   |
| 2    |    |   | Cyclohexene, 1-methyl-4-(1-methy... | 136 | C10H16   | 005989-54-8 | 96   |
| 3    |    |   | Limonene                            | 136 | C10H16   | 000138-86-3 | 91   |
| 4    |    |   | Cyclohexene, 1-methyl-5-(1-methy... | 136 | C10H16   | 013898-73-2 | 81   |
| 5    |    |   | Cyclohexanol, 1-methyl-4-(1-meth... | 196 | C12H20O2 | 010198-23-9 | 80   |



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Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

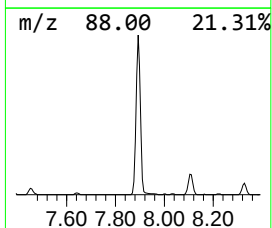
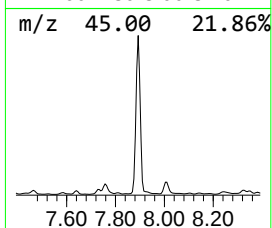
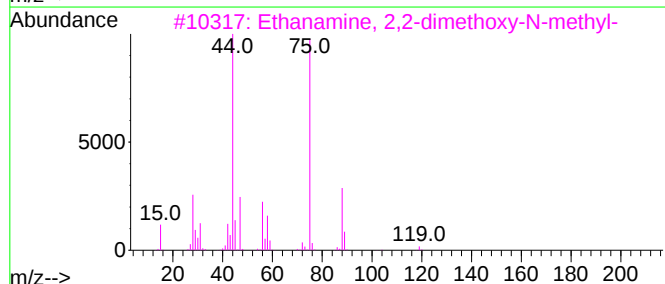
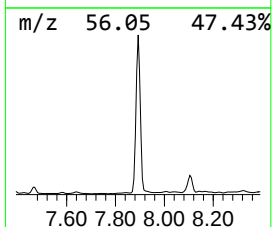
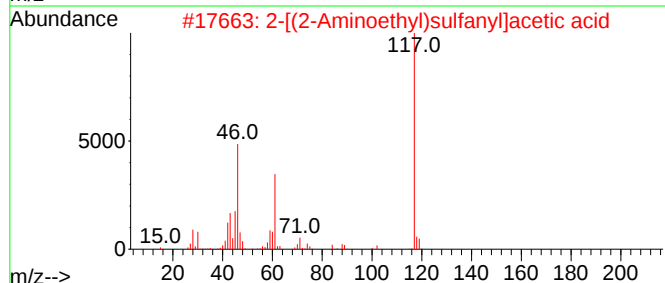
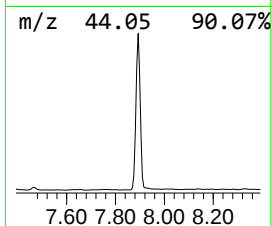
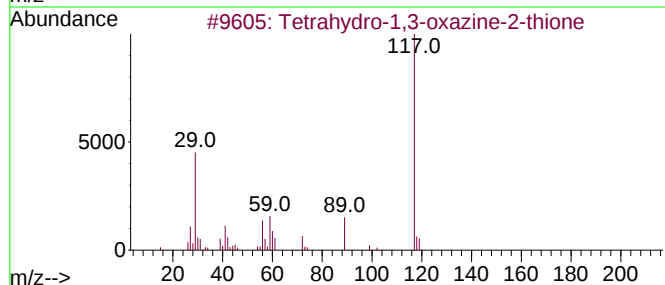
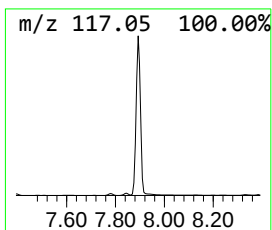
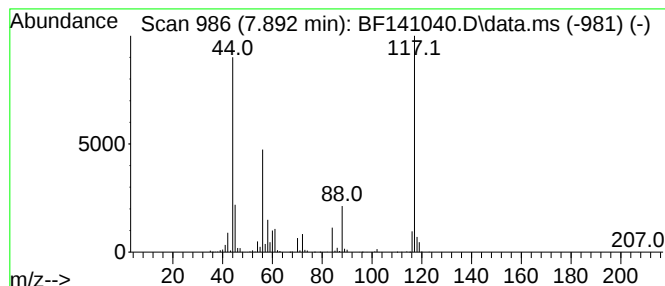
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown7.892 Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 7.892 | 8.84 ng | 815554 | Naphthalene-d8   | 8.104 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Tetrahydro-1,3-oxazine-2-thione     | 117 | C4H7NOS  | 017374-18-4  | 22   |
| 2    |    |   | 2-[(2-Aminoethyl)sulfanyl]acetic... | 135 | C4H9N02S | 003335-52-2  | 12   |
| 3    |    |   | Ethanamine, 2,2-dimethoxy-N-methyl- | 119 | C5H13NO2 | 000122-07-6  | 12   |
| 4    |    |   | 2-Ethylbutyric acid, octyl ester    | 228 | C14H28O2 | 1000340-24-3 | 12   |
| 5    |    |   | Butanoic acid, 3-amino-2-methyl-    | 117 | C5H11NO2 | 032723-74-3  | 12   |



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Misc :  
ALS Vial : 11 Sample Multiplier: 1

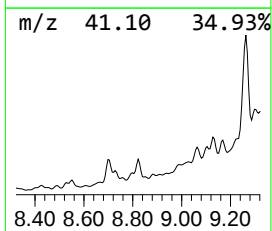
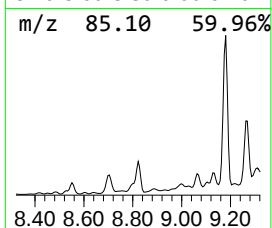
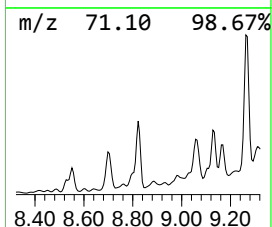
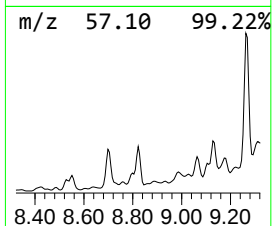
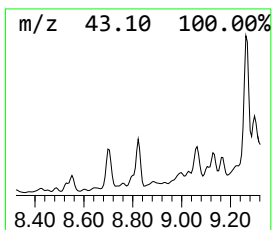
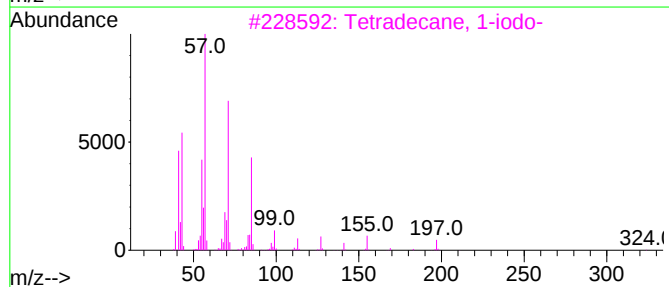
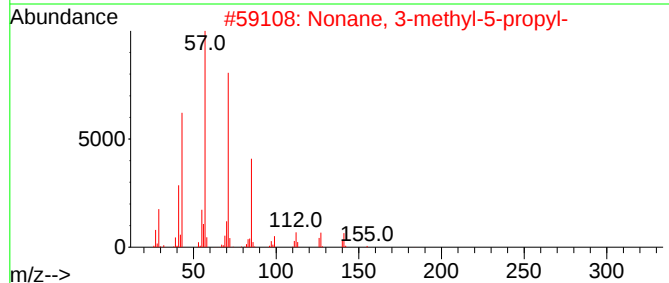
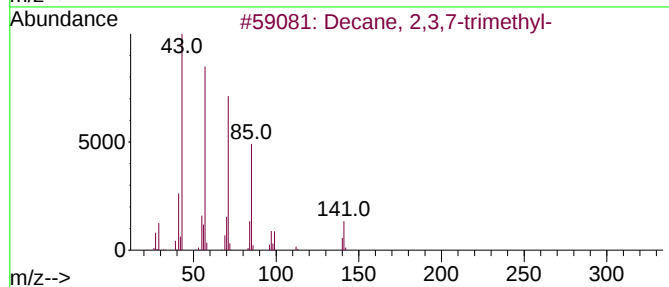
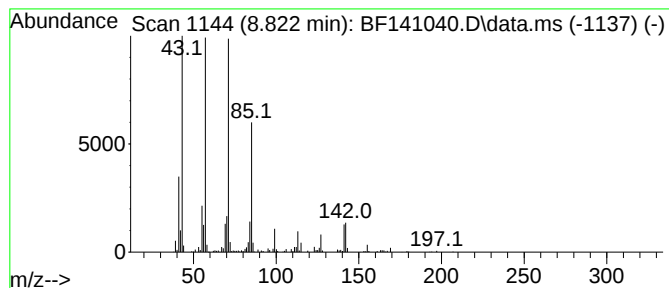
Instrument :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Decane, 2,3,7-trimethyl- Concentration Rank 20

| R.T.      | EstConc                    | Area   | Relative to ISTD | R.T.        |      |
|-----------|----------------------------|--------|------------------|-------------|------|
| 8.822     | 3.43 ng                    | 316558 | Naphthalene-d8   | 8.104       |      |
| Hit# of 5 | Tentative ID               | MW     | MolForm          | CAS#        | Qual |
| 1         | Decane, 2,3,7-trimethyl-   | 184    | C13H28           | 062238-13-5 | 80   |
| 2         | Nonane, 3-methyl-5-propyl- | 184    | C13H28           | 031081-18-2 | 72   |
| 3         | Tetradecane, 1-iodo-       | 324    | C14H29I          | 019218-94-1 | 72   |
| 4         | Dodecane, 1-iodo-          | 296    | C12H25I          | 004292-19-7 | 72   |
| 5         | Dodecane, 4-methyl-        | 184    | C13H28           | 006117-97-1 | 72   |



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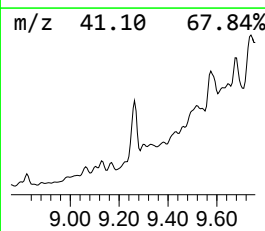
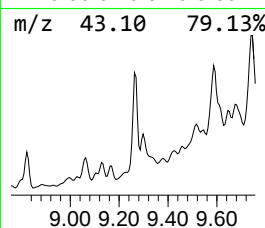
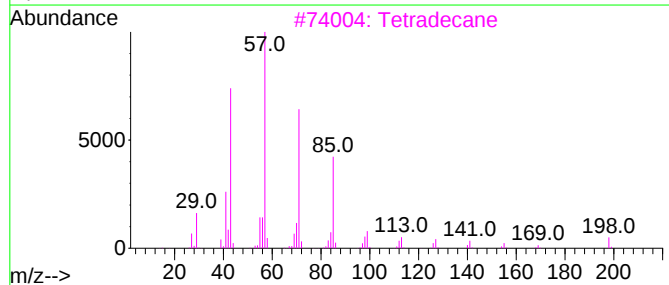
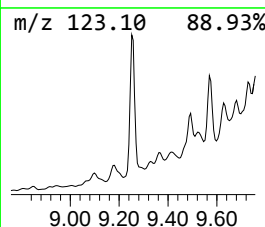
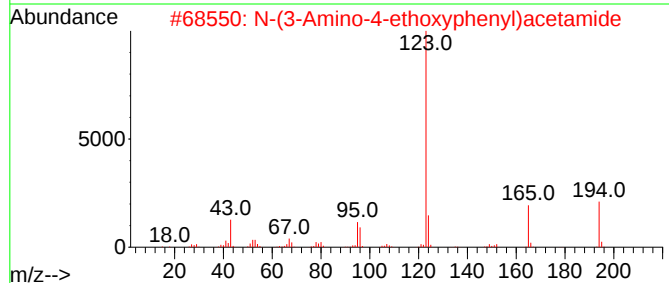
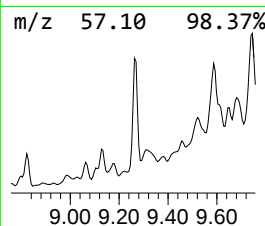
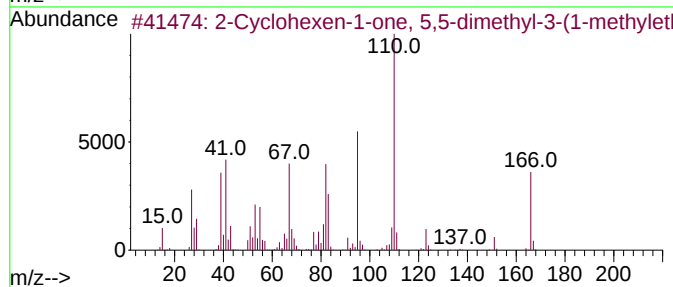
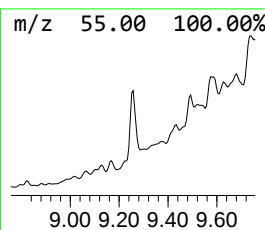
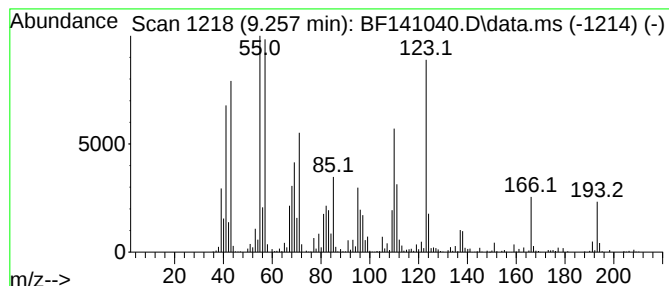
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 unknown9.257 Concentration Rank 11

| R.T.      | EstConc                             | Area    | Relative to ISTD |            | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|------------|-------------|------|
| 9.257     | 10.08 ng                            | 1698930 | Acenaphthene-d10 |            | 9.863       |      |
| Hit# of 5 | Tentative ID                        |         | MW               | MolForm    | CAS#        | Qual |
| 1         | 2-Cyclohexen-1-one, 5,5-dimethyl... |         | 166              | C11H18O    | 028017-79-0 | 35   |
| 2         | N-(3-Amino-4-ethoxyphenyl)acetamide |         | 194              | C10H14N2O2 | 017026-81-2 | 30   |
| 3         | Tetradecane                         |         | 198              | C14H30     | 000629-59-4 | 25   |
| 4         | 1,3-Benzenediol, 4,5-dimethyl-      |         | 138              | C8H10O2    | 000527-55-9 | 18   |
| 5         | 1,4-Hexadiene, 2,3,4,5-tetramethyl- |         | 138              | C10H18     | 051504-54-2 | 18   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

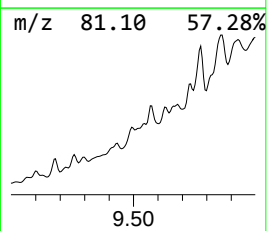
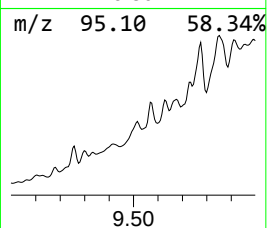
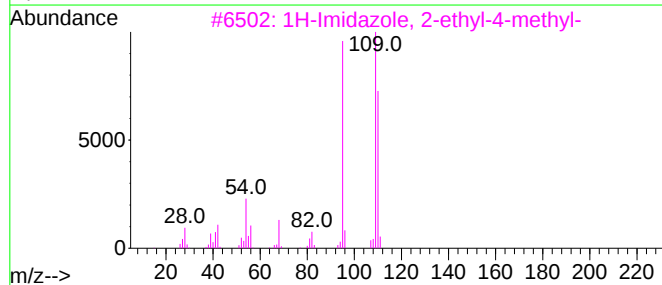
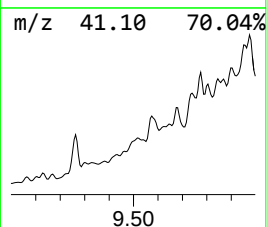
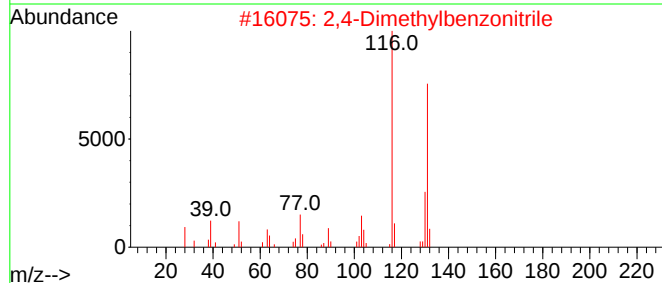
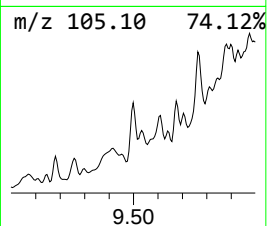
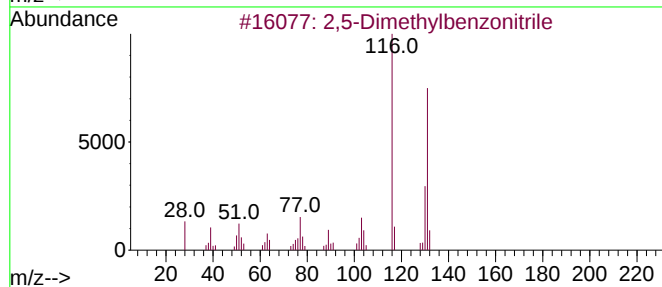
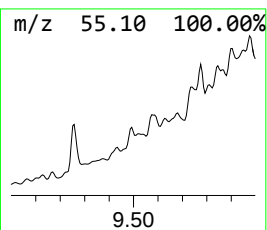
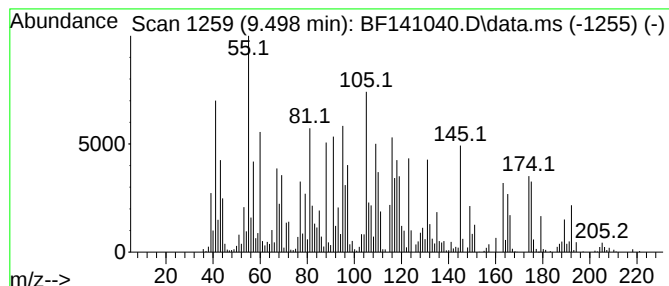
Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 unknown9.498 Concentration Rank 17

| R.T.      | EstConc                          | Area   | Relative to ISTD |          |              | R.T.  |
|-----------|----------------------------------|--------|------------------|----------|--------------|-------|
| 9.498     | 4.50 ng                          | 758186 | Acenaphthene-d10 |          |              | 9.863 |
| Hit# of 5 | Tentative ID                     |        | MW               | MolForm  | CAS#         | Qual  |
| 1         | 2,5-Dimethylbenzonitrile         |        | 131              | C9H9N    | 013730-09-1  | 11    |
| 2         | 2,4-Dimethylbenzonitrile         |        | 131              | C9H9N    | 021789-36-6  | 11    |
| 3         | 1H-Imidazole, 2-ethyl-4-methyl-  |        | 110              | C6H10N2  | 000931-36-2  | 10    |
| 4         | 1,1-Diethoxyhex-trans-4-en-2-yne |        | 168              | C10H16O2 | 1000250-51-9 | 10    |
| 5         | 2,6-Dimethylbenzonitrile         |        | 131              | C9H9N    | 006575-13-9  | 10    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

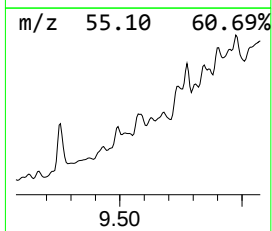
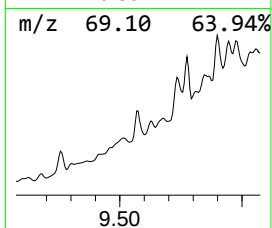
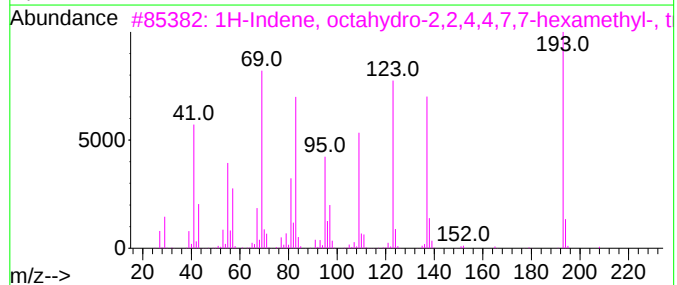
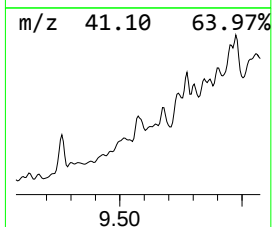
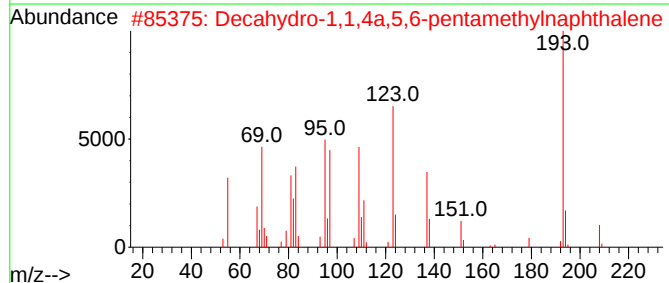
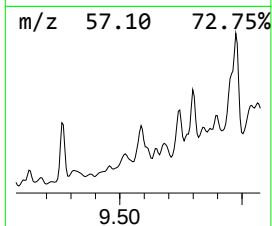
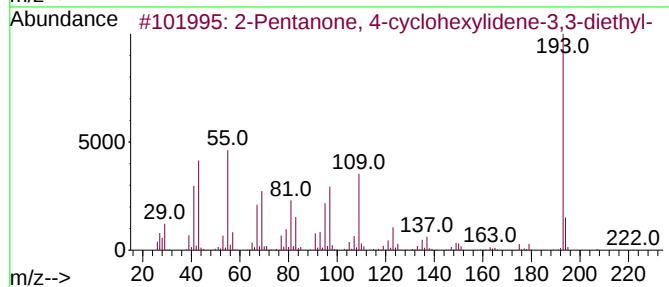
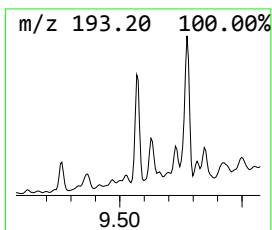
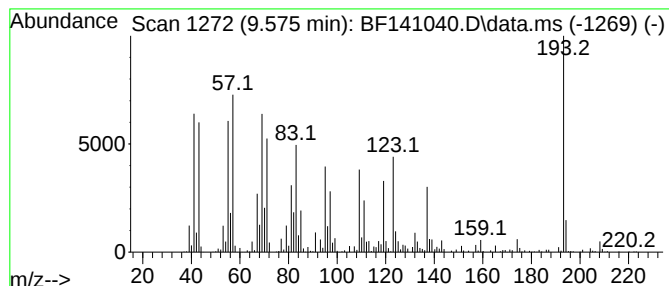
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 2-Pentanone, 4-cyclohexylidene... Concentration Rank 15

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 9.575     | 8.43 ng                             | 1420970 | Acenaphthene-d10 | 9.863        |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | 2-Pentanone, 4-cyclohexylidene-3... | 222     | C15H26O          | 313253-65-5  | 53   |
| 2         | Decahydro-1,1,4a,5,6-pentamethyl... | 208     | C15H28           | 080655-44-3  | 43   |
| 3         | 1H-Indene, octahydro-2,2,4,4,7,7... | 208     | C15H28           | 054832-83-6  | 40   |
| 4         | 1-(p-Fluorophenyl)-4-piperidone     | 193     | C11H12FNO        | 1000238-56-7 | 38   |
| 5         | 2H-Spiro[1-benzofuran-3,2'-[1,3]... | 193     | C10H11NO3        | 1000387-33-4 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

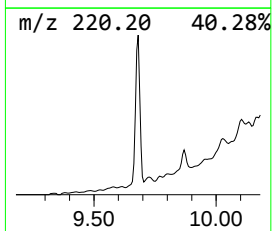
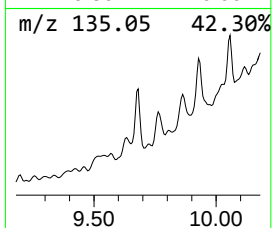
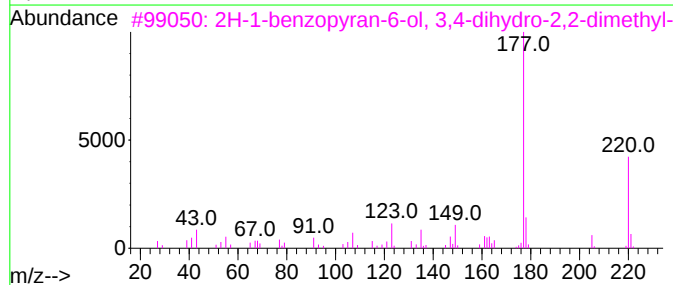
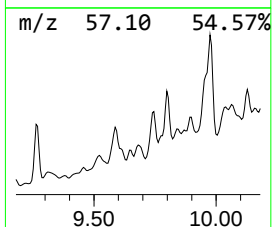
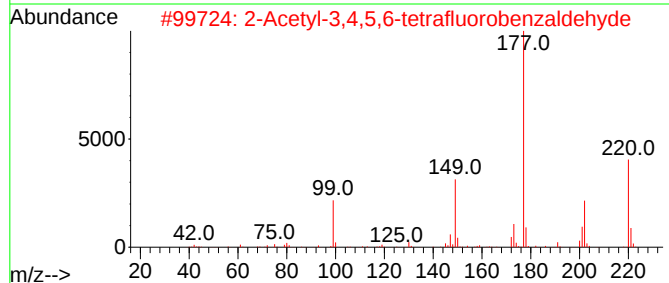
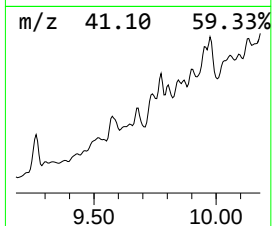
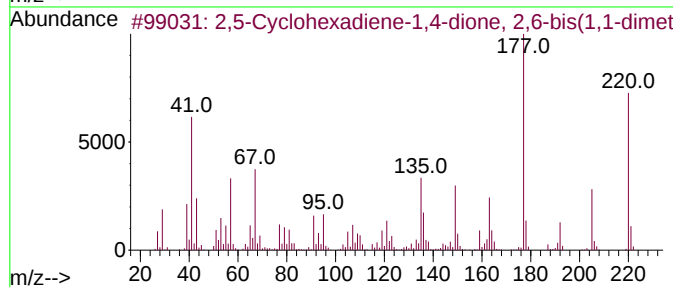
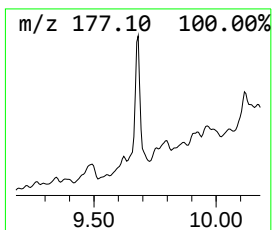
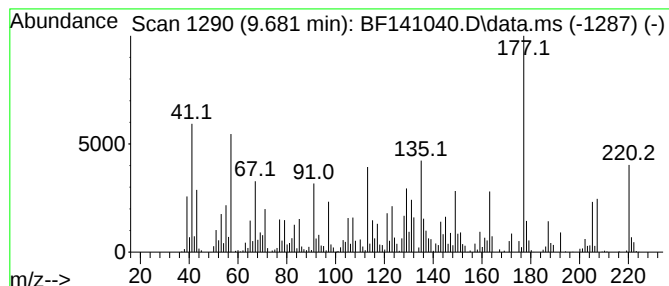
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 2,5-Cyclohexadiene-1,4-dione... Concentration Rank 18

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 9.681     | 4.21 ng                             | 710217 | Acenaphthene-d10 | 9.863        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 2,5-Cyclohexadiene-1,4-dione, 2,... | 220    | C14H2002         | 000719-22-2  | 93   |
| 2         | 2-Acetyl-3,4,5,6-tetrafluorobenz... | 220    | C9H4F402         | 1000461-12-2 | 41   |
| 3         | 2H-1-benzopyran-6-ol, 3,4-dihydr... | 220    | C14H2002         | 1000396-25-4 | 38   |
| 4         | 2-Butanone, 3-(4-tert-butylpheno... | 220    | C14H2002         | 160875-28-5  | 38   |
| 5         | Thiocoumarin, 4,4,6,8-tetramethyl-  | 220    | C13H160S         | 1000128-90-3 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

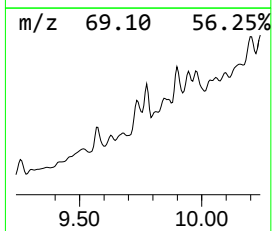
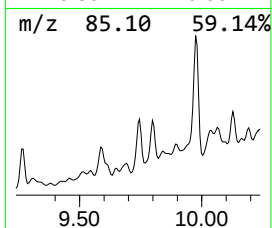
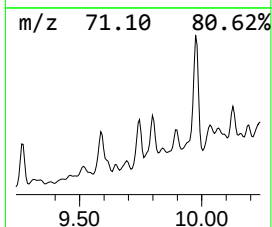
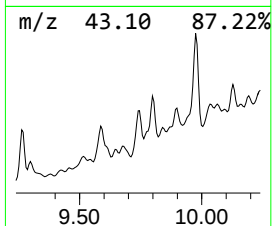
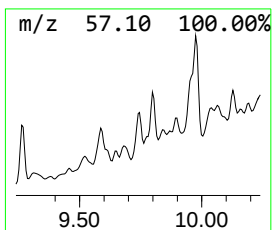
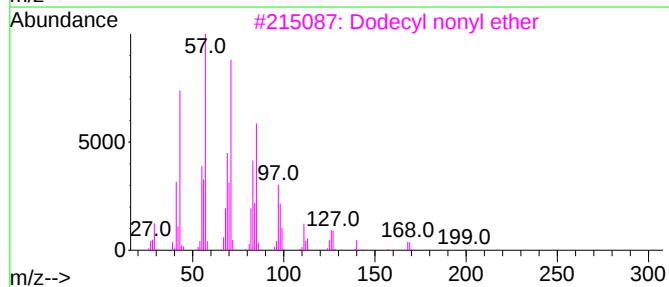
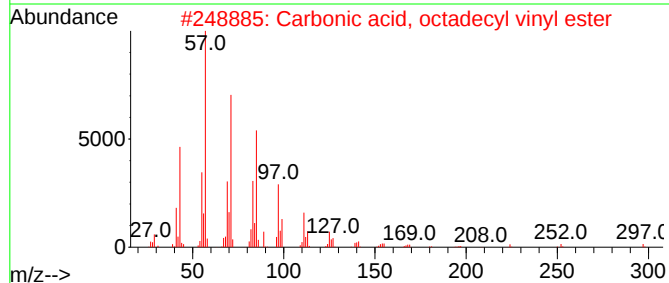
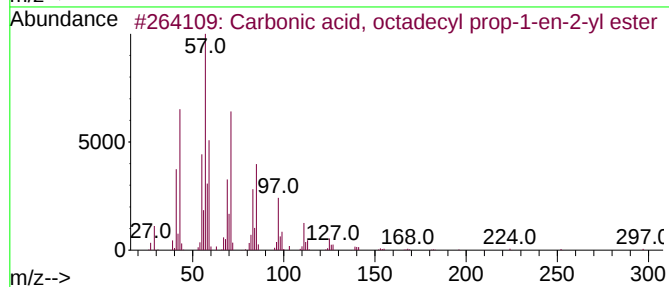
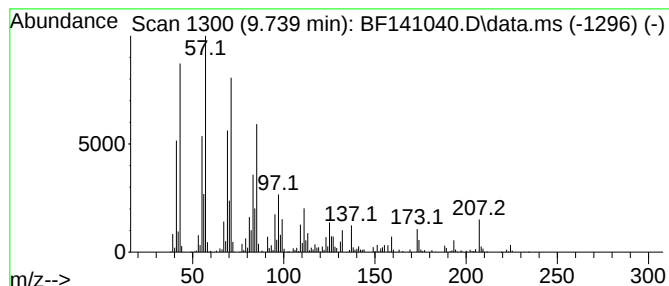
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Carbonic acid, octadecyl pr... Concentration Rank 13

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 9.739     | 9.03 ng                             | 1521570 | Acenaphthene-d10 | 9.863        |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Carbonic acid, octadecyl prop-1-... | 354     | C22H42O3         | 1000383-11-5 | 70   |
| 2         | Carbonic acid, octadecyl vinyl e... | 340     | C21H40O3         | 1000382-54-4 | 68   |
| 3         | Dodecyl nonyl ether                 | 312     | C21H44O          | 1000406-37-5 | 68   |
| 4         | Carbonic acid, eicosyl vinyl ester  | 368     | C23H44O3         | 1000382-54-3 | 62   |
| 5         | Carbonic acid, hexadecyl prop-1-... | 326     | C20H38O3         | 1000382-90-3 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

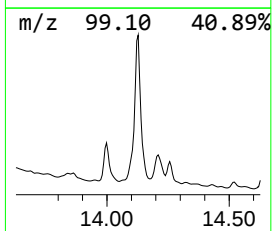
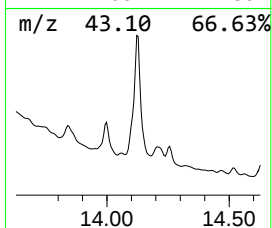
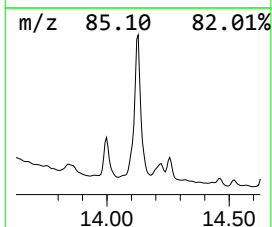
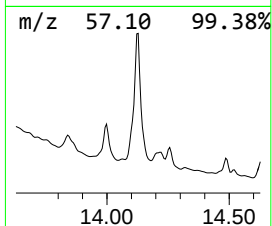
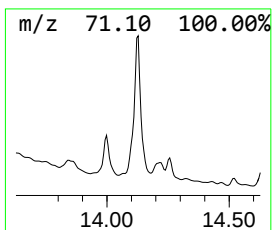
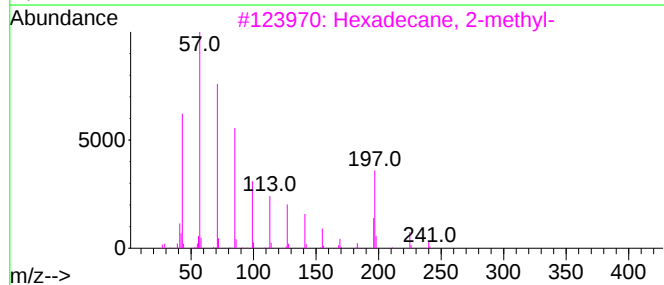
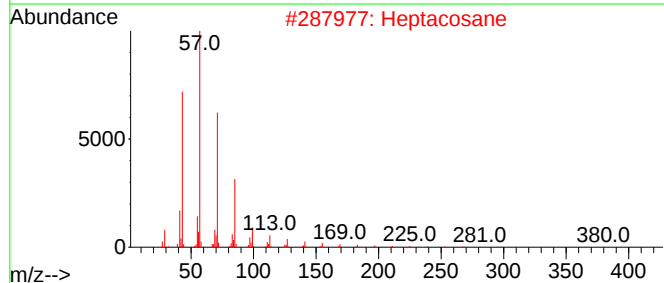
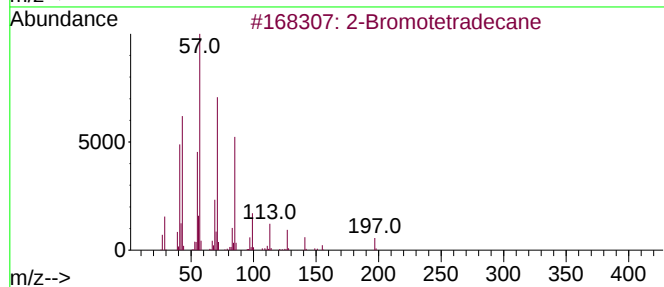
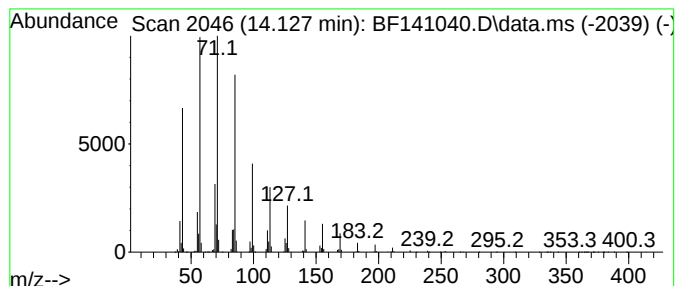
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 2-Bromotetradecane Concentration Rank 5

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 14.127 | 40.11 ng | 5744370 | Chrysene-d12     | 13.992 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm  | CAS#         | Qual |
|------|------|------------------------|-----|----------|--------------|------|
| 1    |      | 2-Bromotetradecane     | 276 | C14H29Br | 074036-95-6  | 83   |
| 2    |      | Heptacosane            | 380 | C27H56   | 000593-49-7  | 72   |
| 3    |      | Hexadecane, 2-methyl-  | 240 | C17H36   | 001560-92-5  | 72   |
| 4    |      | Dotriacontane, 1-iodo- | 576 | C32H65I  | 1000406-32-4 | 53   |
| 5    |      | 9-Methylheneicosane    | 310 | C22H46   | 070475-51-3  | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
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Misc :  
ALS Vial : 11 Sample Multiplier: 1

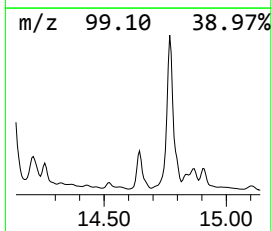
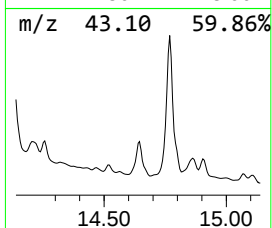
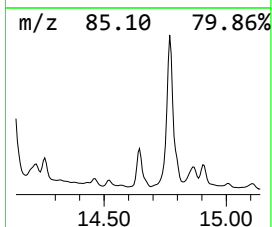
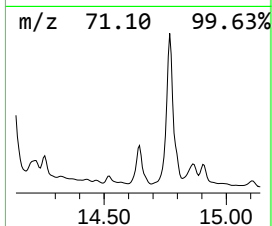
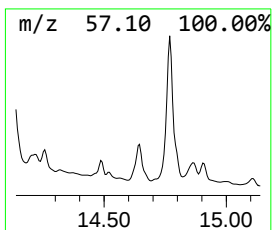
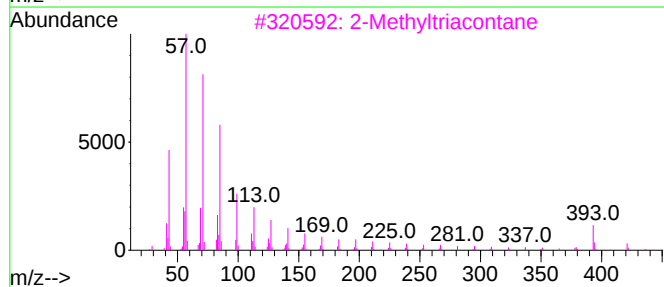
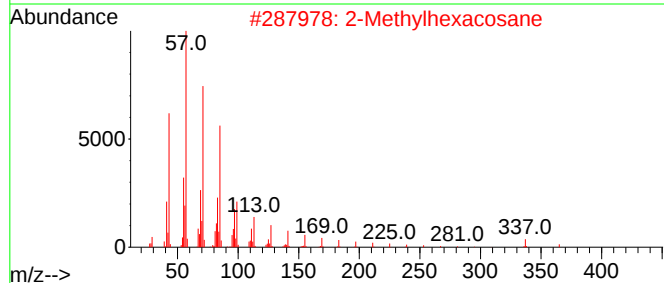
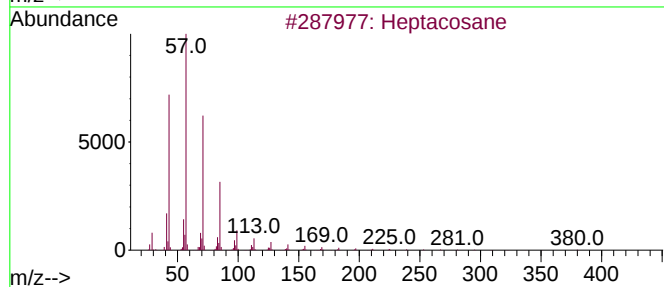
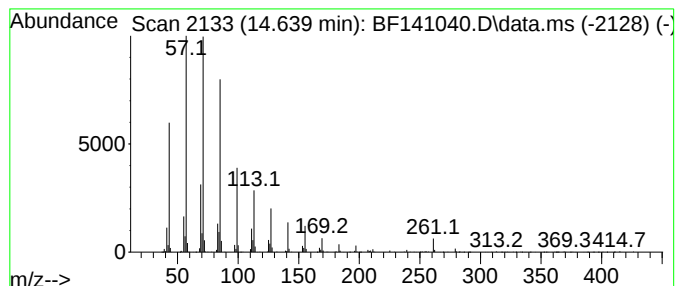
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ClientSampleId :  
MOO-24-00395-96

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Heptacosane Concentration Rank 9

| R.T.      | EstConc                      | Area    | Relative to ISTD |         | R.T.         |      |
|-----------|------------------------------|---------|------------------|---------|--------------|------|
| 14.639    | 13.45 ng                     | 1926330 | Chrysene-d12     |         | 13.992       |      |
| Hit# of 5 | Tentative ID                 |         | MW               | MolForm | CAS#         | Qual |
| 1         | Heptacosane                  |         | 380              | C27H56  | 000593-49-7  | 83   |
| 2         | 2-Methylhexacosane           |         | 380              | C27H56  | 001561-02-0  | 72   |
| 3         | 2-Methyltriacontane          |         | 437              | C31H64  | 001560-72-1  | 72   |
| 4         | Octacosane, 1-iodo-          |         | 520              | C28H57I | 1000406-32-2 | 64   |
| 5         | Dodecane, 2-methyl-6-propyl- |         | 226              | C16H34  | 055045-08-4  | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

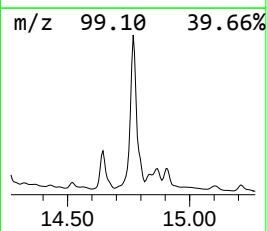
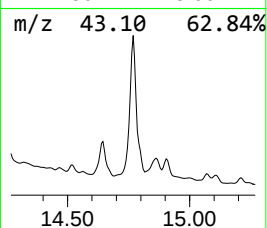
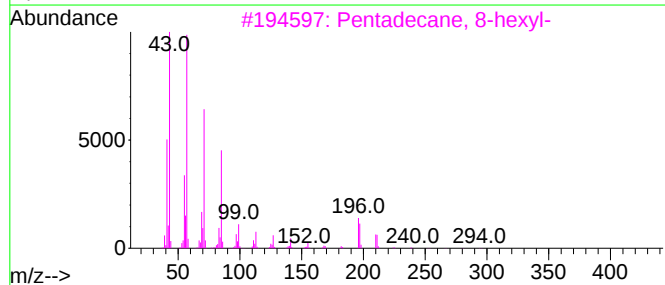
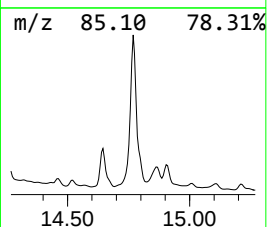
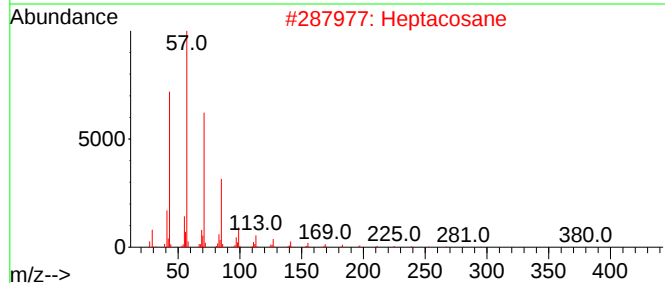
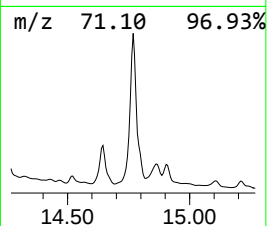
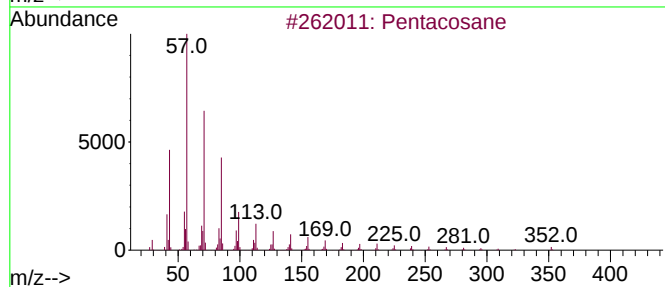
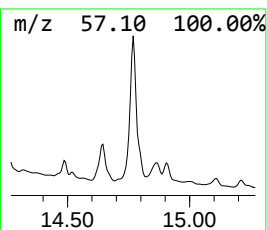
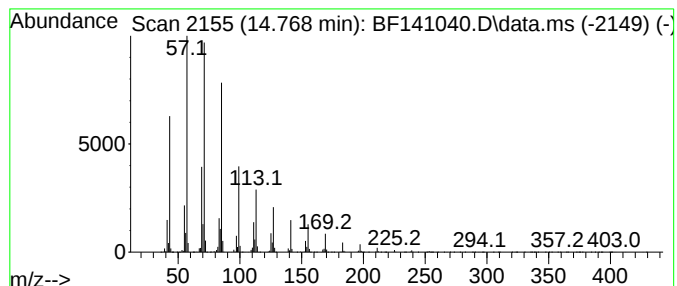
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 Pentacosane Concentration Rank 2

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 14.769 | 83.41 ng | 5836670 | Perylene-d12     | 15.445 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Pentacosane           | 352 | C25H52  | 000629-99-2 | 90   |
| 2    |    |   | Heptacosane           | 380 | C27H56  | 000593-49-7 | 86   |
| 3    |    |   | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 81   |
| 4    |    |   | Tetradecane, 1-iodo-  | 324 | C14H29I | 019218-94-1 | 74   |
| 5    |    |   | Decane, 3,8-dimethyl- | 170 | C12H26  | 017312-55-9 | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

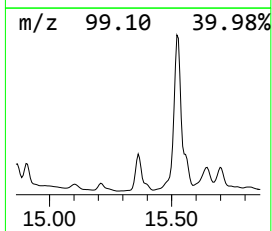
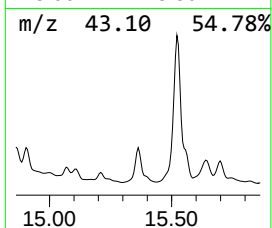
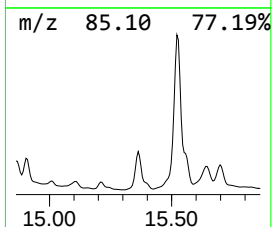
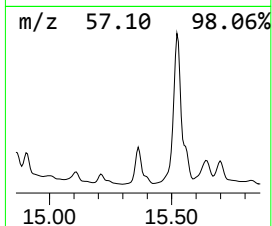
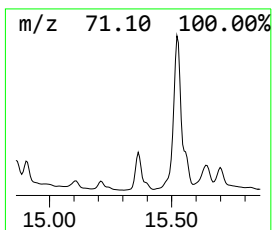
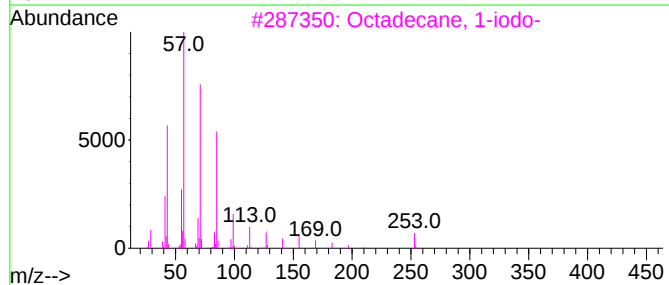
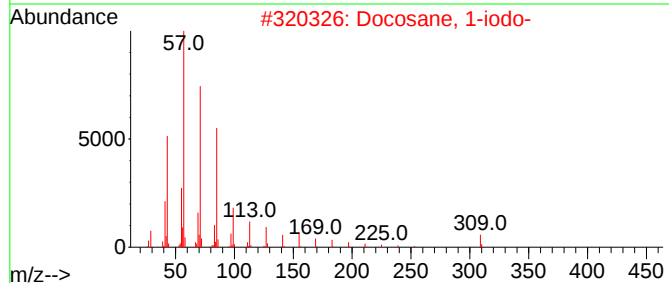
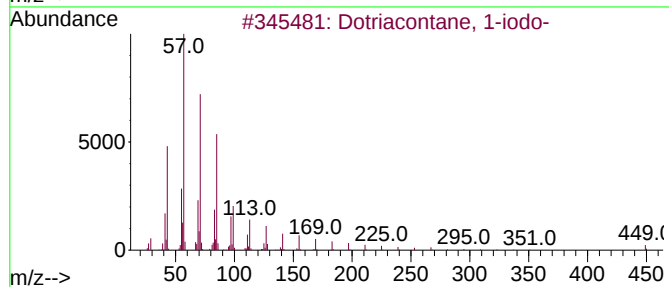
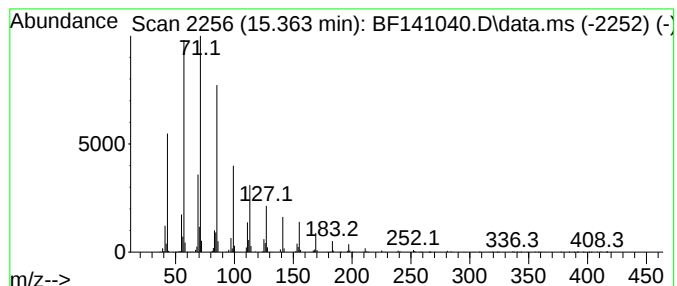
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Dotriacontane, 1-iodo- Concentration Rank 8

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 15.363 | 13.90 ng | 972643 | Perylene-d12     | 15.445 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#         | Qual |
|------|----|---|------------------------|-----|---------|--------------|------|
| 1    |    |   | Dotriacontane, 1-iodo- | 576 | C32H65I | 1000406-32-4 | 72   |
| 2    |    |   | Docosane, 1-iodo-      | 436 | C22H45I | 1000406-31-9 | 59   |
| 3    |    |   | Octadecane, 1-iodo-    | 380 | C18H37I | 000629-93-6  | 58   |
| 4    |    |   | Dodecane, 4-methyl-    | 184 | C13H28  | 006117-97-1  | 58   |
| 5    |    |   | Eicosane, 1-iodo-      | 408 | C20H41I | 1000406-31-8 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

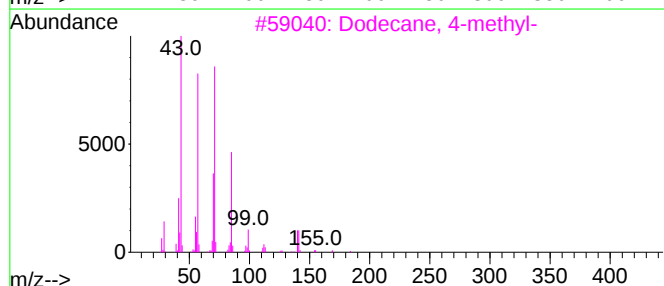
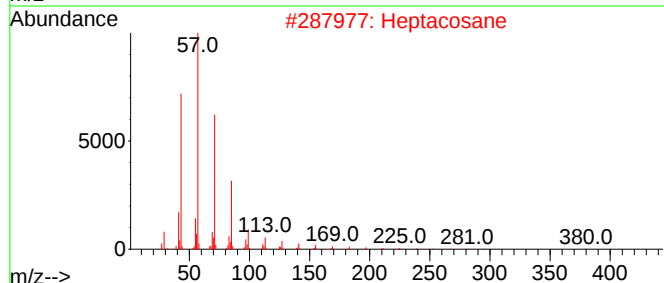
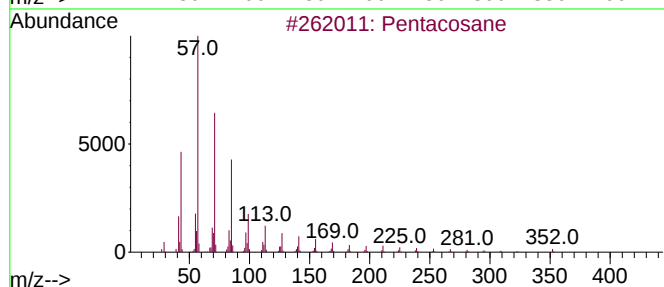
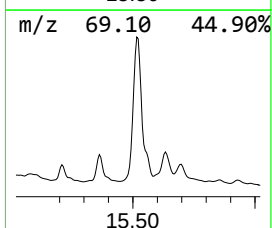
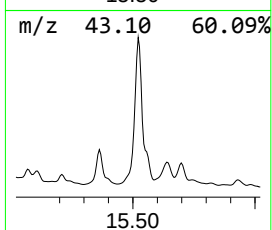
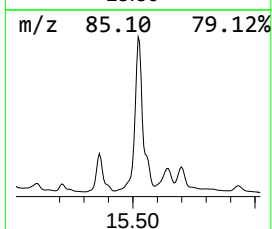
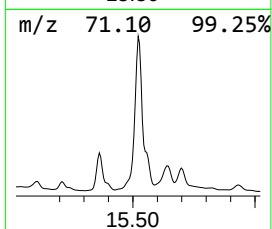
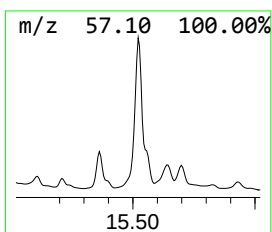
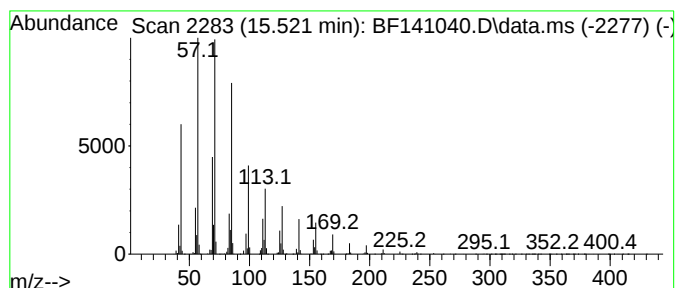
Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.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 14 Dodecane, 4-methyl- Concentration Rank 1

| R.T.      | EstConc                      | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|------------------------------|---------|------------------|---------|-------------|------|
| 15.521    | 85.13 ng                     | 5957350 | Perylene-d12     |         | 15.445      |      |
| Hit# of 5 | Tentative ID                 |         | MW               | MolForm | CAS#        | Qual |
| 1         | Pentacosane                  |         | 352              | C25H52  | 000629-99-2 | 90   |
| 2         | Heptacosane                  |         | 380              | C27H56  | 000593-49-7 | 90   |
| 3         | Dodecane, 4-methyl-          |         | 184              | C13H28  | 006117-97-1 | 53   |
| 4         | Dodecane, 2-methyl-6-propyl- |         | 226              | C16H34  | 055045-08-4 | 50   |
| 5         | Octane, 5-ethyl-2-methyl-    |         | 156              | C11H24  | 062016-18-6 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

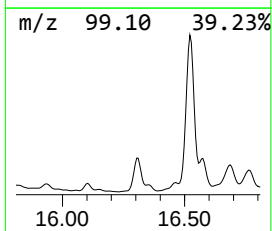
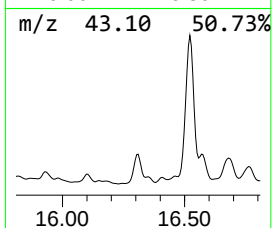
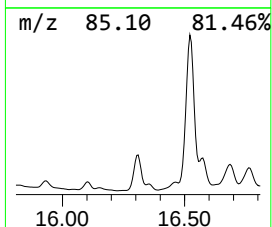
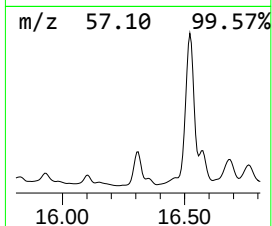
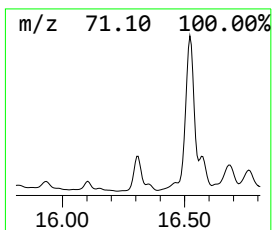
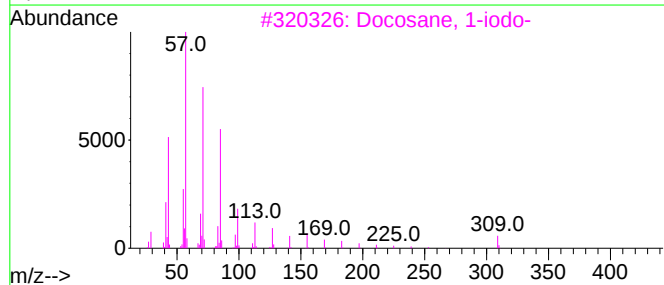
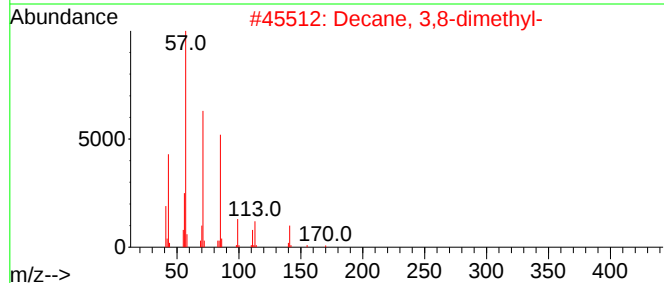
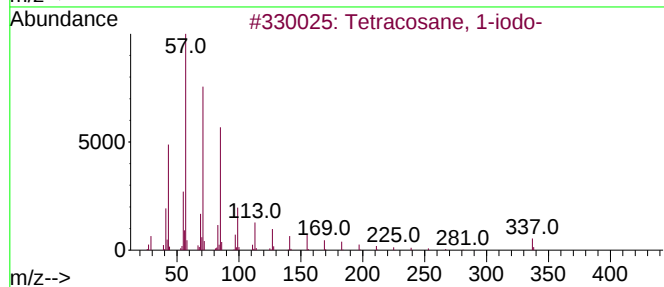
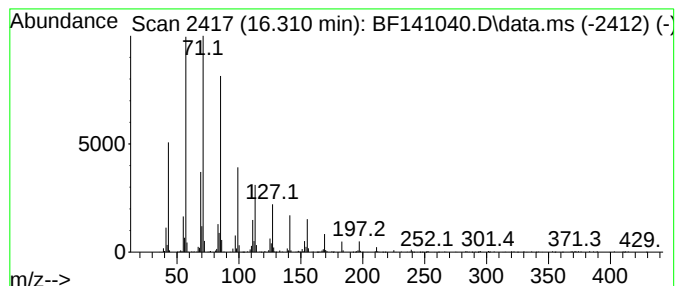
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Tetracosane, 1-iodo- Concentration Rank 12

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 16.310 | 9.78 ng | 684622 | Perylene-d12     | 15.445 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#         | Qual |
|------|----|---|-----------------------|-----|---------|--------------|------|
| 1    |    |   | Tetracosane, 1-iodo-  | 464 | C24H49I | 1000406-32-0 | 72   |
| 2    |    |   | Decane, 3,8-dimethyl- | 170 | C12H26  | 017312-55-9  | 68   |
| 3    |    |   | Docosane, 1-iodo-     | 436 | C22H45I | 1000406-31-9 | 64   |
| 4    |    |   | Eicosane, 1-iodo-     | 408 | C20H41I | 1000406-31-8 | 59   |
| 5    |    |   | Decane, 3-methyl-     | 156 | C11H24  | 013151-34-3  | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

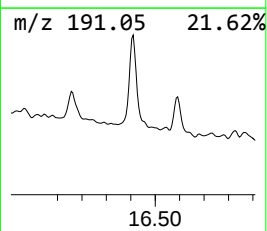
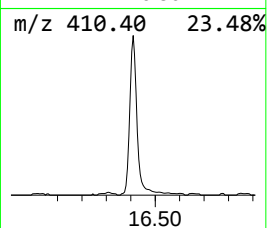
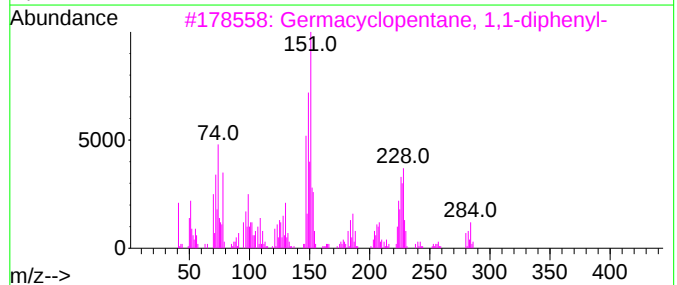
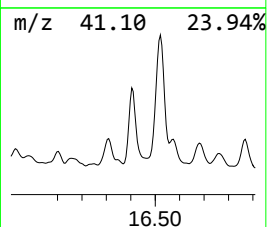
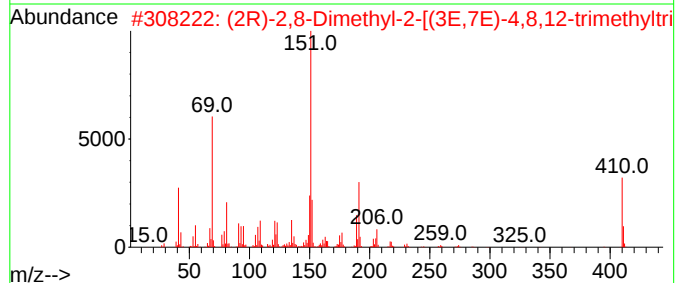
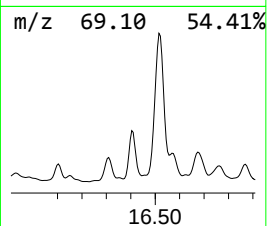
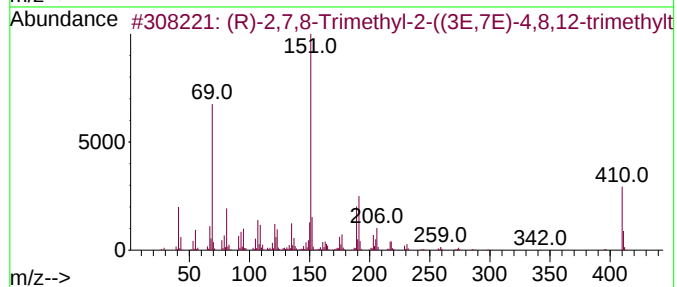
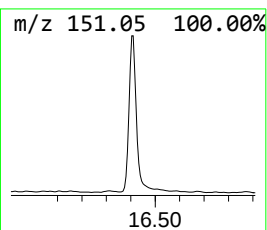
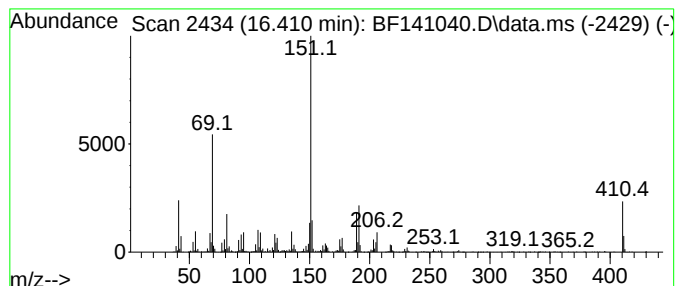
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 (R)-2,7,8-Trimethyl-2-((3E,... Concentration Rank 10

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 16.410 | 12.23 ng | 855807 | Perylene-d12     | 15.445 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | (R)-2,7,8-Trimethyl-2-((3E,7E)-4... | 410 | C28H42O2     | 014101-61-2  | 99      |      |      |
| 2    | (2R)-2,8-Dimethyl-2-[(3E,7E)-4,8... | 410 | C28H42O2     | 1000503-47-0 | 99      |      |      |
| 3    | Germacypentane, 1,1-diphenyl-       | 284 | C16H18Ge     | 004514-06-1  | 87      |      |      |
| 4    | .gamma.-Tocopherol                  | 416 | C28H48O2     | 007616-22-0  | 58      |      |      |
| 5    | Benzenepropanenitrile, 3,4-dimet... | 191 | C11H13NO2    | 049621-56-9  | 47      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

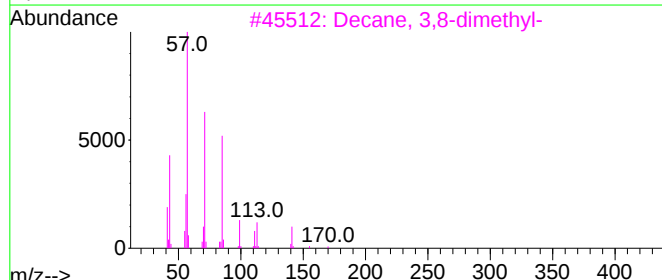
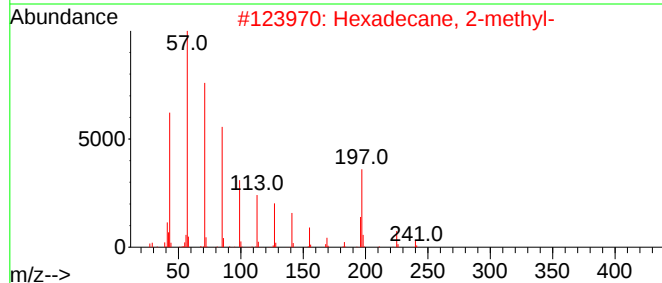
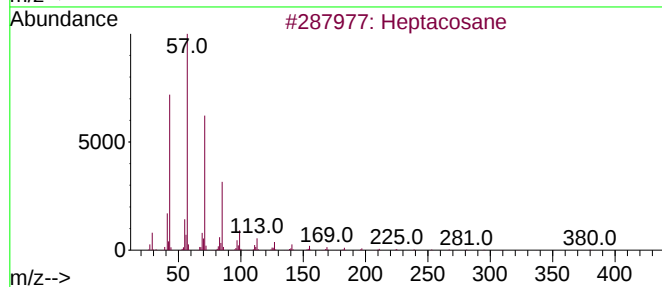
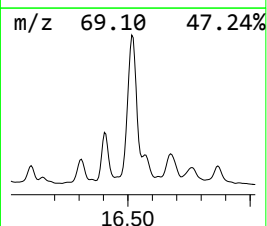
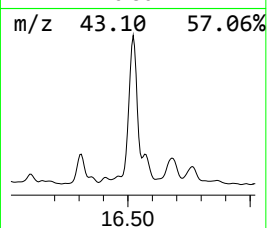
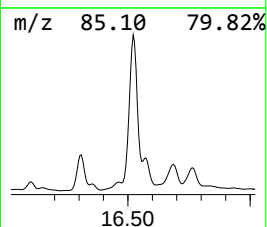
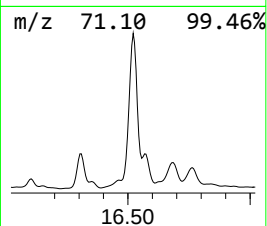
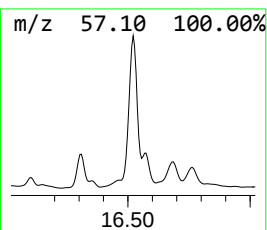
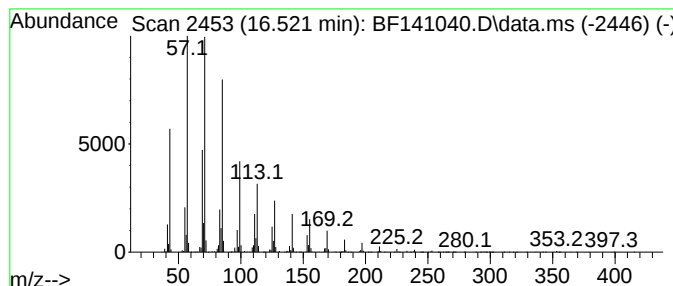
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 17 Hexadecane, 2-methyl- Concentration Rank 3

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 16.521 | 67.83 ng | 4746830 | Perylene-d12     | 15.445 |

| Hit# | of 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------------|-----|---------|-------------|------|
| 1    |      | Heptacosane               | 380 | C27H56  | 000593-49-7 | 90   |
| 2    |      | Hexadecane, 2-methyl-     | 240 | C17H36  | 001560-92-5 | 86   |
| 3    |      | Decane, 3,8-dimethyl-     | 170 | C12H26  | 017312-55-9 | 68   |
| 4    |      | Tridecane, 7-propyl-      | 226 | C16H34  | 055045-09-5 | 53   |
| 5    |      | Octane, 5-ethyl-2-methyl- | 156 | C11H24  | 062016-18-6 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

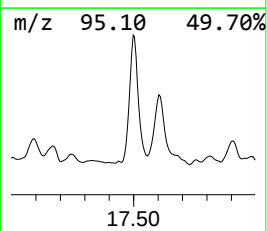
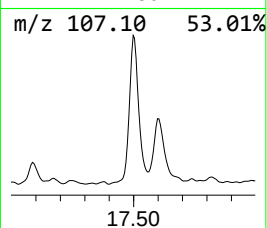
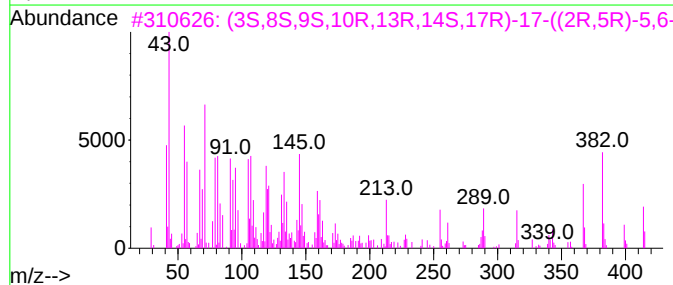
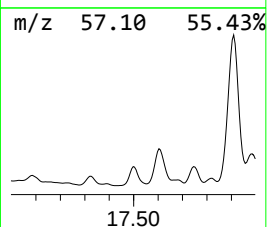
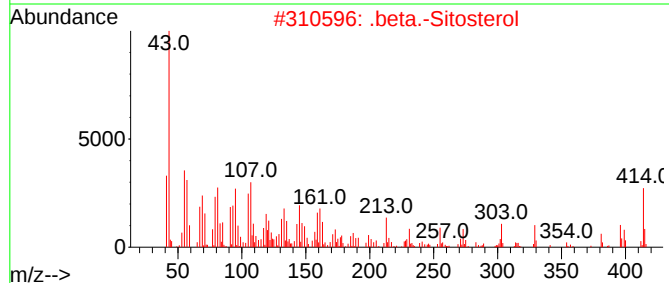
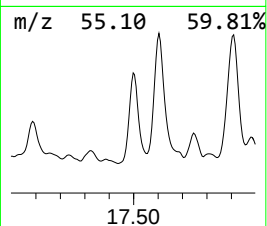
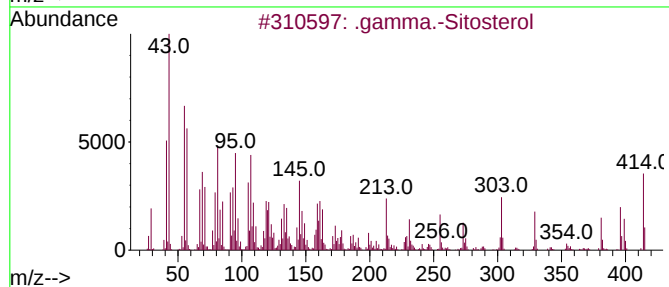
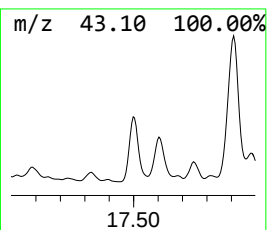
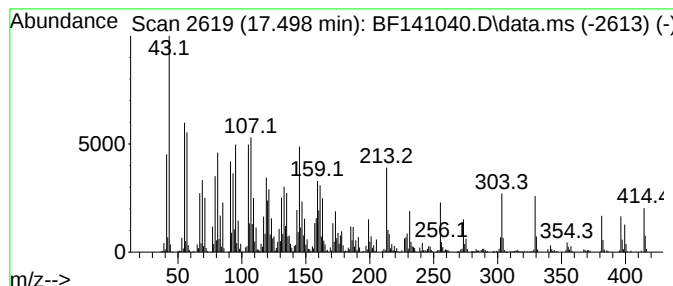
Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.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 18 .gamma.-Sitosterol Concentration Rank 7

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 17.498    | 29.00 ng                            | 2029580 | Perylene-d12     | 15.445      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | .gamma.-Sitosterol                  | 414     | C29H50O          | 000083-47-6 | 99   |
| 2         | .beta.-Sitosterol                   | 414     | C29H50O          | 000083-46-5 | 94   |
| 3         | (3S,8S,9S,10R,13R,14S,17R)-17-((... | 414     | C29H50O          | 029365-29-5 | 55   |
| 4         | Campesterol                         | 400     | C28H48O          | 000474-62-4 | 49   |
| 5         | Ergost-7-en-3-ol, (3.beta.)-        | 400     | C28H48O          | 026047-31-4 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

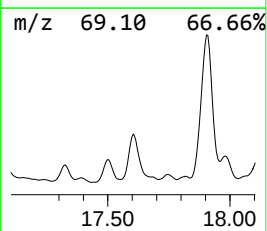
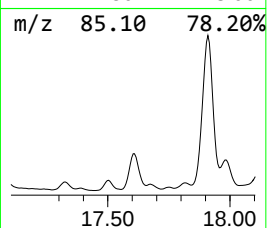
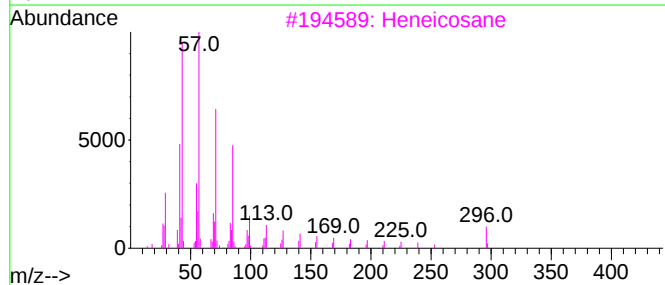
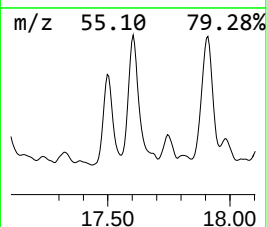
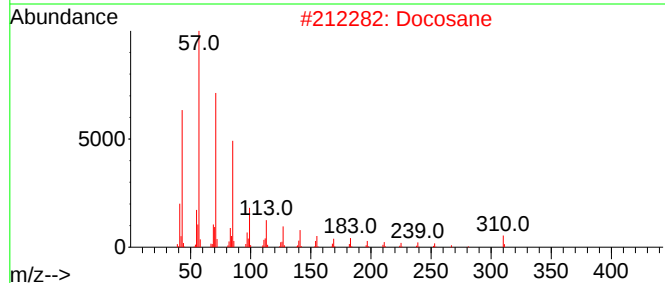
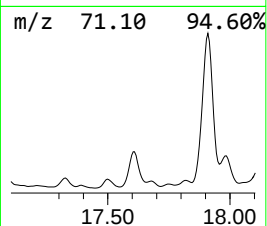
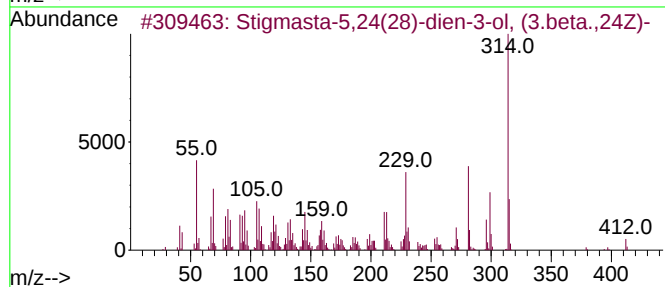
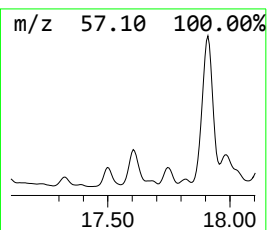
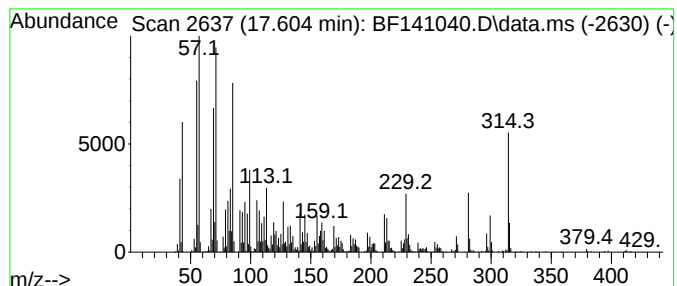
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 Stigmasta-5,24(28)-dien-3-o... Concentration Rank 6

| R.T.      | EstConc                             | Area    | Relative to ISTD |         | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|---------|--------------|------|
| 17.604    | 31.85 ng                            | 2228590 | Perylene-d12     |         | 15.445       |      |
| Hit# of 5 | Tentative ID                        |         | MW               | MolForm | CAS#         | Qual |
| 1         | Stigmasta-5,24(28)-dien-3-ol, (3... |         | 412              | C29H48O | 000481-14-1  | 93   |
| 2         | Docosane                            |         | 310              | C22H46  | 000629-97-0  | 91   |
| 3         | Heneicosane                         |         | 296              | C21H44  | 000629-94-7  | 83   |
| 4         | Heptadecane, 2-methyl-              |         | 254              | C18H38  | 001560-89-0  | 83   |
| 5         | Octacosane, 1-iodo-                 |         | 520              | C28H57I | 1000406-32-2 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

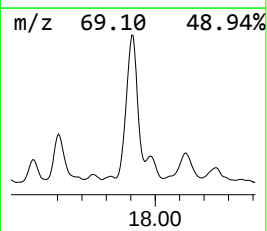
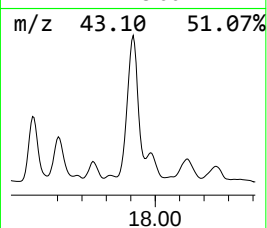
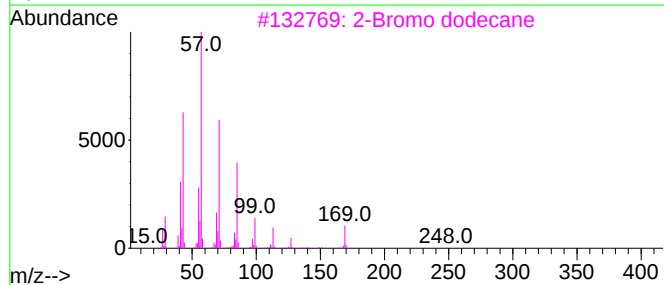
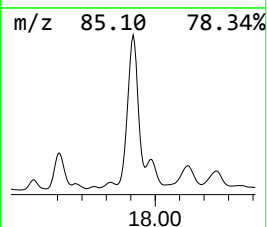
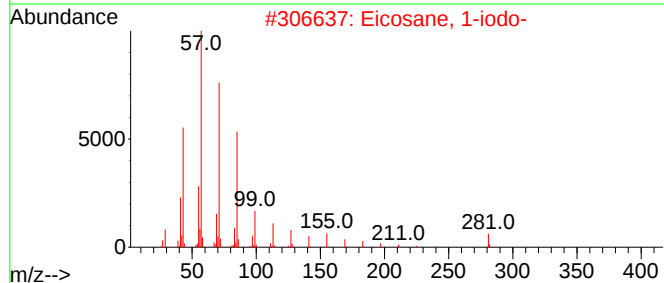
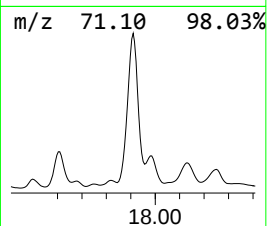
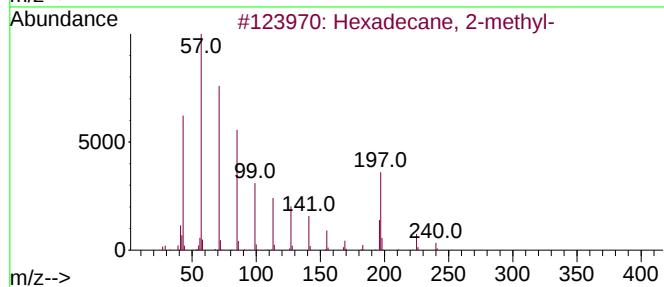
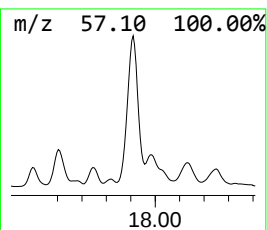
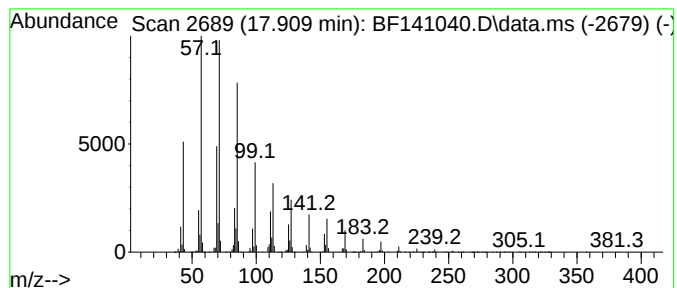
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 20 Eicosane, 1-iodo- Concentration Rank 4

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 17.910 | 57.16 ng | 3999800 | Perylene-d12     | 15.445 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm  | CAS#         | Qual |
|------|------|------------------------------|-----|----------|--------------|------|
| 1    |      | Hexadecane, 2-methyl-        | 240 | C17H36   | 001560-92-5  | 86   |
| 2    |      | Eicosane, 1-iodo-            | 408 | C20H41I  | 1000406-31-8 | 53   |
| 3    |      | 2-Bromo dodecane             | 248 | C12H25Br | 013187-99-0  | 53   |
| 4    |      | Dodecane, 2-methyl-6-propyl- | 226 | C16H34   | 055045-08-4  | 50   |
| 5    |      | Octane, 5-ethyl-2-methyl-    | 156 | C11H24   | 062016-18-6  | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF123024\  
Data File : BF141040.D  
Acq On : 30 Dec 2024 19:25  
Operator : RC/JU  
Sample : P5386-03  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
MOO-24-00395-96

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF122624.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 |
| Pentanal           | 2.399  | 5.1     | ng    | 357977   | 1                     | 6.828  | 1400970 | 20.0 |
| D-Limonene         | 6.945  | 4.1     | ng    | 289766   | 1                     | 6.828  | 1400970 | 20.0 |
| unknown7.892       | 7.892  | 8.8     | ng    | 815554   | 2                     | 8.104  | 1845790 | 20.0 |
| Decane, 2,3,7-t... | 8.822  | 3.4     | ng    | 316558   | 2                     | 8.104  | 1845790 | 20.0 |
| unknown9.257       | 9.257  | 10.1    | ng    | 1698930  | 3                     | 9.863  | 3371000 | 20.0 |
| unknown9.498       | 9.498  | 4.5     | ng    | 758186   | 3                     | 9.863  | 3371000 | 20.0 |
| 2-Pentanone, 4-... | 9.575  | 8.4     | ng    | 1420970  | 3                     | 9.863  | 3371000 | 20.0 |
| 2,5-Cyclohexadi... | 9.681  | 4.2     | ng    | 710217   | 3                     | 9.863  | 3371000 | 20.0 |
| Carbonic acid, ... | 9.739  | 9.0     | ng    | 1521570  | 3                     | 9.863  | 3371000 | 20.0 |
| 2-Bromotetradecane | 14.127 | 40.1    | ng    | 5744370  | 5                     | 13.992 | 2864000 | 20.0 |
| Heptacosane        | 14.639 | 13.4    | ng    | 1926330  | 5                     | 13.992 | 2864000 | 20.0 |
| Pentacosane        | 14.769 | 83.4    | ng    | 5836670  | 6                     | 15.445 | 1399530 | 20.0 |
| Dotriacontane, ... | 15.363 | 13.9    | ng    | 972643   | 6                     | 15.445 | 1399530 | 20.0 |
| Dodecane, 4-met... | 15.521 | 85.1    | ng    | 5957350  | 6                     | 15.445 | 1399530 | 20.0 |
| Tetracosane, 1-... | 16.310 | 9.8     | ng    | 684622   | 6                     | 15.445 | 1399530 | 20.0 |
| (R)-2,7,8-Trime... | 16.410 | 12.2    | ng    | 855807   | 6                     | 15.445 | 1399530 | 20.0 |
| Hexadecane, 2-m... | 16.521 | 67.8    | ng    | 4746830  | 6                     | 15.445 | 1399530 | 20.0 |
| .gamma.-Sitosterol | 17.498 | 29.0    | ng    | 2029580  | 6                     | 15.445 | 1399530 | 20.0 |
| Stigmasta-5,24(... | 17.604 | 31.9    | ng    | 2228590  | 6                     | 15.445 | 1399530 | 20.0 |
| Eicosane, 1-iodo-  | 17.910 | 57.2    | ng    | 3999800  | 6                     | 15.445 | 1399530 | 20.0 |