

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
 Data File : BF142767.D  
 Acq On : 12 Jun 2025 19:56  
 Operator : RC/JU  
 Sample : Q2195-01 5X  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OK-01-060325

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-BF061125.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142767.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         | 5.098       | 488           | 494         | 506          | rBV      | 193335         | 271146        | 37.37%          | 4.139%        |
| 2         | 5.504       | 559           | 563         | 576          | rBV      | 81998          | 112468        | 15.50%          | 1.717%        |
| 3         | 6.516       | 730           | 735         | 749          | rBV      | 68745          | 94863         | 13.08%          | 1.448%        |
| 4         | 6.892       | 794           | 799         | 805          | rBV      | 194920         | 235974        | 32.53%          | 3.602%        |
| 5         | 7.451       | 890           | 894         | 905          | rBV      | 41095          | 55247         | 7.62%           | 0.843%        |
| 6         | 8.175       | 1011          | 1017        | 1025         | rBV      | 199843         | 254849        | 35.13%          | 3.890%        |
| 7         | 9.251       | 1195          | 1200        | 1210         | rBV      | 99882          | 129527        | 17.85%          | 1.977%        |
| 8         | 9.586       | 1253          | 1257        | 1260         | rVV      | 6995           | 10168         | 1.40%           | 0.155%        |
| 9         | 9.645       | 1264          | 1267        | 1273         | rVV4     | 5342           | 10712         | 1.48%           | 0.164%        |
| 10        | 9.704       | 1273          | 1277        | 1283         | rVB      | 4986           | 7618          | 1.05%           | 0.116%        |
| 11        | 9.792       | 1288          | 1292        | 1297         | rBV      | 20658          | 26624         | 3.67%           | 0.406%        |
| 12        | 9.851       | 1297          | 1302        | 1307         | rVB3     | 4693           | 9146          | 1.26%           | 0.140%        |
| 13        | 9.933       | 1311          | 1316        | 1321         | rBV      | 242297         | 302932        | 41.76%          | 4.624%        |
| 14        | 10.133      | 1345          | 1350        | 1354         | rBV2     | 9847           | 14814         | 2.04%           | 0.226%        |
| 15        | 10.175      | 1354          | 1357        | 1362         | rVB      | 11986          | 14101         | 1.94%           | 0.215%        |
| 16        | 10.245      | 1365          | 1369        | 1372         | rBV2     | 13009          | 17246         | 2.38%           | 0.263%        |
| 17        | 10.275      | 1372          | 1374        | 1380         | rVB3     | 6578           | 8709          | 1.20%           | 0.133%        |
| 18        | 10.357      | 1380          | 1388        | 1391         | rBV4     | 10683          | 24783         | 3.42%           | 0.378%        |
| 19        | 10.480      | 1405          | 1409        | 1415         | rVB3     | 11003          | 17074         | 2.35%           | 0.261%        |
| 20        | 10.592      | 1425          | 1428        | 1431         | rVB3     | 5906           | 7274          | 1.00%           | 0.111%        |
| 21        | 10.639      | 1431          | 1436        | 1443         | rVB3     | 12661          | 21604         | 2.98%           | 0.330%        |
| 22        | 10.722      | 1444          | 1450        | 1455         | rBV      | 72368          | 96425         | 13.29%          | 1.472%        |
| 23        | 10.839      | 1465          | 1470        | 1476         | rVB4     | 21541          | 36640         | 5.05%           | 0.559%        |
| 24        | 10.922      | 1480          | 1484        | 1487         | rBV      | 7945           | 11269         | 1.55%           | 0.172%        |
| 25        | 11.027      | 1497          | 1502        | 1505         | rBV2     | 13780          | 23996         | 3.31%           | 0.366%        |
| 26        | 11.057      | 1505          | 1507        | 1510         | rVB2     | 9674           | 9805          | 1.35%           | 0.150%        |
| 27        | 11.110      | 1514          | 1516        | 1520         | rVB      | 8742           | 9343          | 1.29%           | 0.143%        |
| 28        | 11.163      | 1520          | 1525        | 1526         | rBV4     | 8687           | 14042         | 1.94%           | 0.214%        |
| 29        | 11.227      | 1532          | 1536        | 1539         | rBV3     | 9737           | 12952         | 1.79%           | 0.198%        |
| 30        | 11.274      | 1540          | 1544        | 1547         | rVV4     | 12420          | 20210         | 2.79%           | 0.309%        |
| 31        | 11.316      | 1547          | 1551        | 1556         | rVB2     | 16790          | 28039         | 3.86%           | 0.428%        |
| 32        | 11.421      | 1564          | 1569        | 1572         | rBV      | 297563         | 422475        | 58.23%          | 6.449%        |
| 33        | 11.498      | 1578          | 1582        | 1585         | rBV      | 21603          | 25850         | 3.56%           | 0.395%        |
| 34        | 11.533      | 1585          | 1588        | 1591         | rVB3     | 10439          | 13239         | 1.82%           | 0.202%        |
| 35        | 11.657      | 1605          | 1609        | 1614         | rBV5     | 12127          | 24385         | 3.36%           | 0.372%        |
| 36        | 11.710      | 1615          | 1618        | 1622         | rVV2     | 16544          | 26622         | 3.67%           | 0.406%        |

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 Operator : RC/JU  
 Sample : Q2195-01 5X  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OK-01-060325

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-BF061125.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |         |         |
|----|--------|------|------|------|------|--------|--------|---------|---------|
| 37 | 11.751 | 1622 | 1625 | 1631 | rVB  | 25714  | 35663  | 4.92%   | 0.544%  |
| 38 | 11.833 | 1636 | 1639 | 1643 | rVB  | 13965  | 18588  | 2.56%   | 0.284%  |
| 39 | 11.880 | 1643 | 1647 | 1650 | rBV3 | 8984   | 13276  | 1.83%   | 0.203%  |
| 40 | 11.927 | 1650 | 1655 | 1657 | rBV  | 68896  | 104404 | 14.39%  | 1.594%  |
| 41 | 11.951 | 1657 | 1659 | 1664 | rVB  | 92441  | 103893 | 14.32%  | 1.586%  |
| 42 | 11.998 | 1664 | 1667 | 1669 | rBV2 | 21471  | 26582  | 3.66%   | 0.406%  |
| 43 | 12.033 | 1669 | 1673 | 1683 | rVB2 | 110904 | 229576 | 31.64%  | 3.505%  |
| 44 | 12.116 | 1683 | 1687 | 1691 | rBV3 | 16804  | 27955  | 3.85%   | 0.427%  |
| 45 | 12.157 | 1691 | 1694 | 1697 | rVB2 | 16333  | 18449  | 2.54%   | 0.282%  |
| 46 | 12.221 | 1697 | 1705 | 1707 | rBV2 | 46731  | 79391  | 10.94%  | 1.212%  |
| 47 | 12.321 | 1715 | 1722 | 1724 | rBV6 | 10336  | 24462  | 3.37%   | 0.373%  |
| 48 | 12.368 | 1724 | 1730 | 1732 | rBV2 | 26388  | 48130  | 6.63%   | 0.735%  |
| 49 | 12.398 | 1732 | 1735 | 1738 | rBV  | 27796  | 32891  | 4.53%   | 0.502%  |
| 50 | 12.474 | 1743 | 1748 | 1751 | rBV  | 107483 | 158034 | 21.78%  | 2.412%  |
| 51 | 12.510 | 1751 | 1754 | 1760 | rVV2 | 58163  | 98517  | 13.58%  | 1.504%  |
| 52 | 12.563 | 1760 | 1763 | 1771 | rVB3 | 49093  | 79642  | 10.98%  | 1.216%  |
| 53 | 12.639 | 1771 | 1776 | 1780 | rBV  | 286442 | 349495 | 48.17%  | 5.335%  |
| 54 | 12.680 | 1780 | 1783 | 1785 | rBV  | 17440  | 22835  | 3.15%   | 0.349%  |
| 55 | 12.792 | 1799 | 1802 | 1805 | rBV  | 18484  | 20056  | 2.76%   | 0.306%  |
| 56 | 12.868 | 1809 | 1815 | 1821 | rBV  | 384254 | 554490 | 76.43%  | 8.465%  |
| 57 | 12.921 | 1821 | 1824 | 1828 | rVB2 | 40274  | 51100  | 7.04%   | 0.780%  |
| 58 | 13.004 | 1834 | 1838 | 1847 | rVB  | 177697 | 284132 | 39.16%  | 4.337%  |
| 59 | 13.086 | 1847 | 1852 | 1858 | rBV4 | 65331  | 127356 | 17.55%  | 1.944%  |
| 60 | 13.139 | 1859 | 1861 | 1863 | rVV2 | 16751  | 17992  | 2.48%   | 0.275%  |
| 61 | 13.192 | 1864 | 1870 | 1874 | rVB2 | 70928  | 144329 | 19.89%  | 2.203%  |
| 62 | 13.263 | 1880 | 1882 | 1885 | rVB  | 29752  | 27578  | 3.80%   | 0.421%  |
| 63 | 13.298 | 1885 | 1888 | 1893 | rBV  | 68052  | 85163  | 11.74%  | 1.300%  |
| 64 | 13.392 | 1901 | 1904 | 1907 | rBV  | 58074  | 68517  | 9.44%   | 1.046%  |
| 65 | 13.427 | 1907 | 1910 | 1913 | rVB  | 46183  | 54190  | 7.47%   | 0.827%  |
| 66 | 14.068 | 2013 | 2019 | 2030 | rBV3 | 334605 | 725475 | 100.00% | 11.075% |
| 67 | 15.127 | 2195 | 2199 | 2206 | rBV  | 66271  | 134454 | 18.53%  | 2.053%  |
| 68 | 15.498 | 2259 | 2262 | 2268 | rVB  | 62631  | 92877  | 12.80%  | 1.418%  |
| 69 | 15.568 | 2269 | 2274 | 2287 | rVB  | 146547 | 268975 | 37.08%  | 4.106%  |

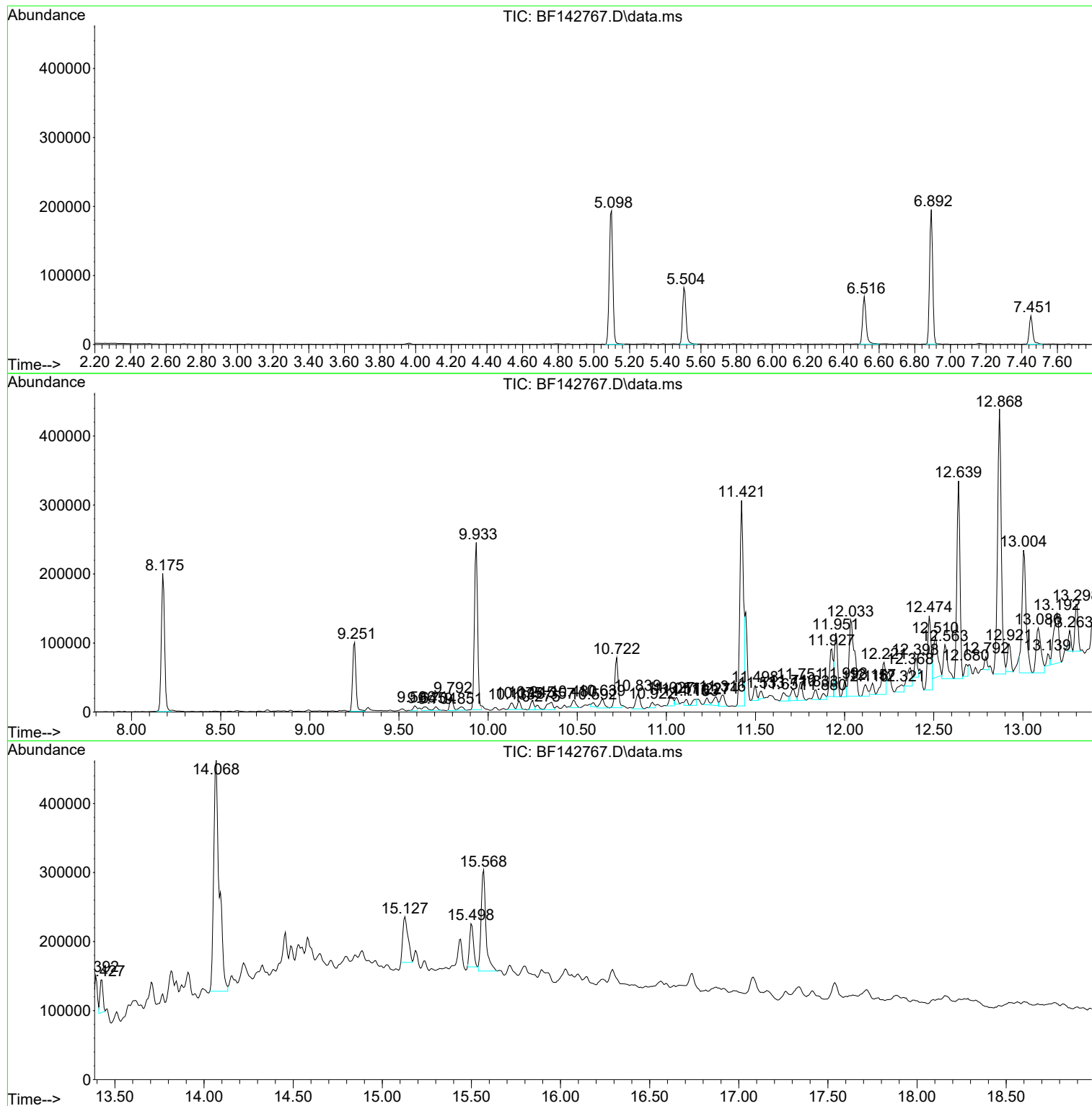
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
Data File : BF142767.D  
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Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF061125.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\BF061225\  
Data File : BF142767.D  
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Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF061125.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

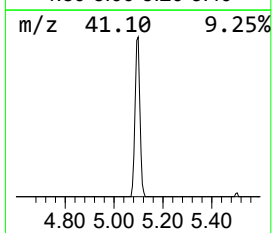
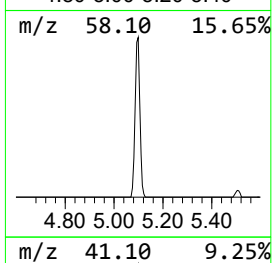
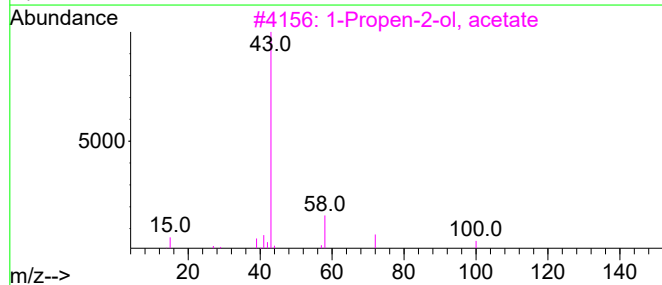
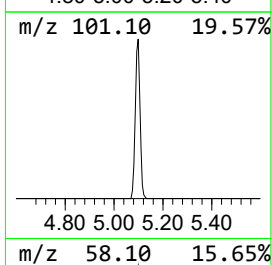
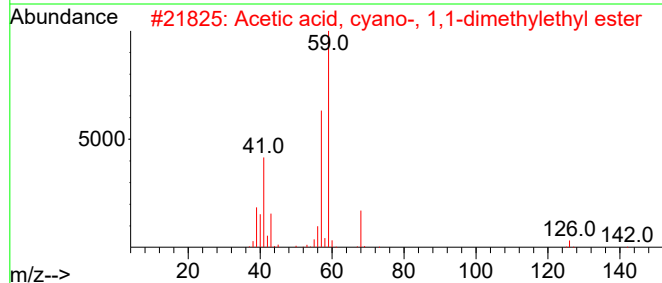
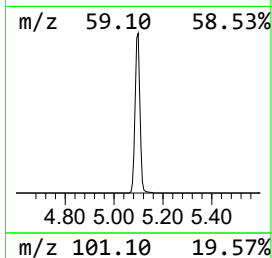
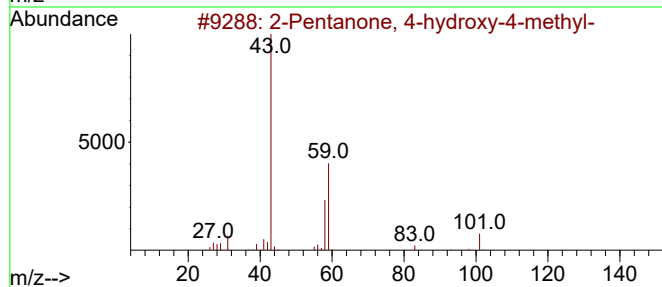
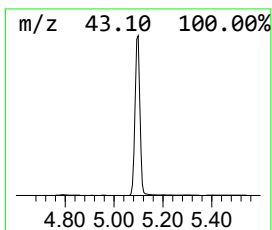
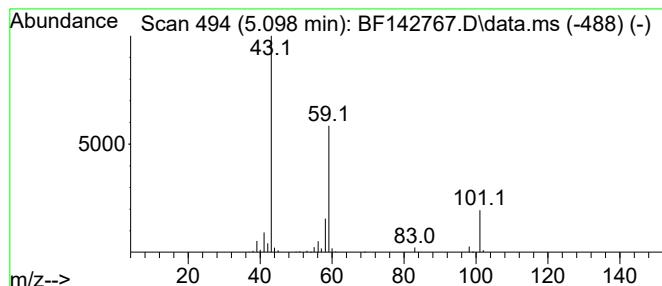
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 5.098 | 22.98 ng | 271146 | 1,4-Dichlorobenzene-d4 | 6.892 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2      | 000123-42-2 | 56      |      |      |
| 2    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2     | 001116-98-9 | 23      |      |      |
| 3    | 1-Propen-2-ol, acetate              | 100 | C5H8O2       | 000108-22-5 | 10      |      |      |
| 4    | Morpholine, 4-methyl-               | 101 | C5H11NO      | 000109-02-4 | 9       |      |      |
| 5    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2      | 000057-71-6 | 9       |      |      |



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Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

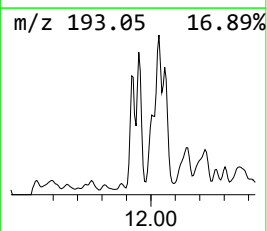
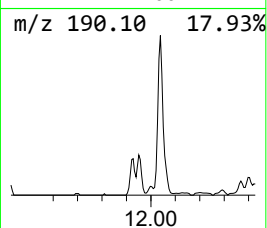
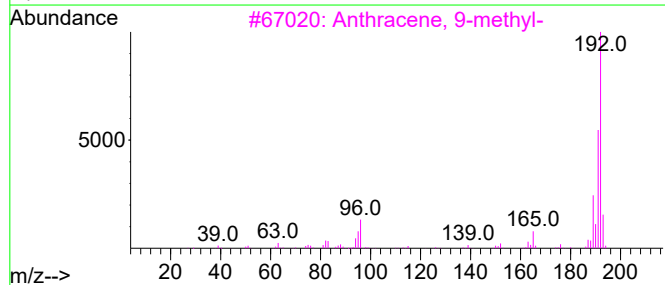
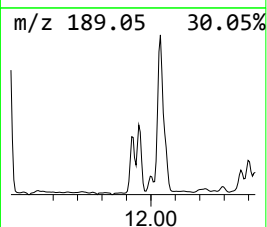
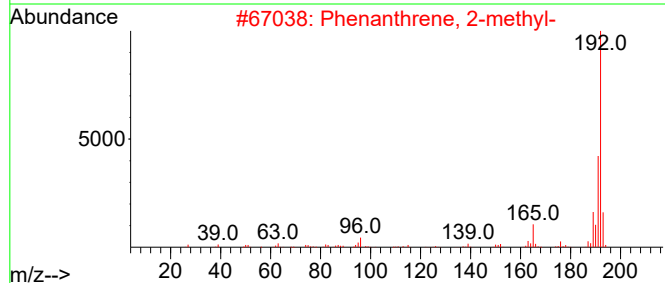
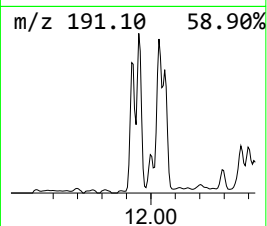
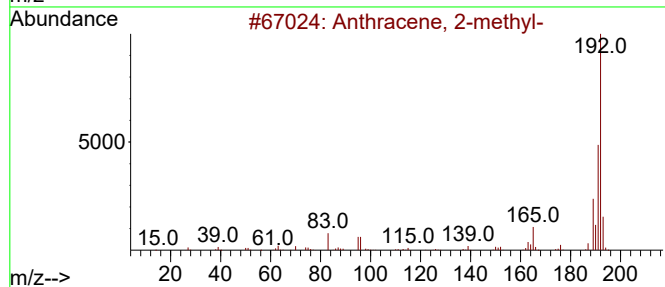
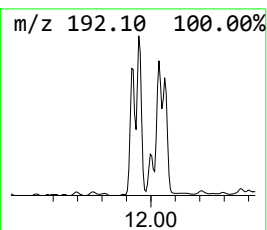
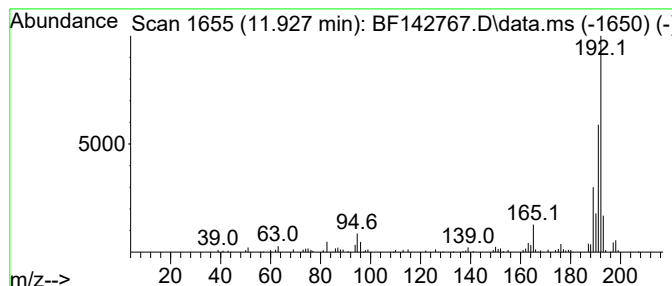
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF061125.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 Anthracene, 2-methyl- Concentration Rank 4

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.927 | 4.94 ng | 104404 | Phenanthrene-d10 | 11.422 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 93   |
| 2    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 91   |
| 3    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 90   |
| 4    |    |   | 1H-Indene, 2-phenyl-    | 192 | C15H12  | 004505-48-0 | 90   |
| 5    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 90   |



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Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

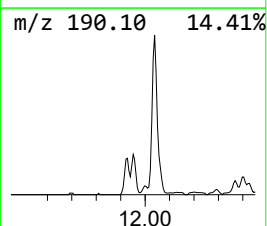
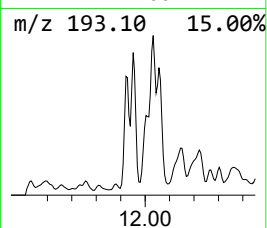
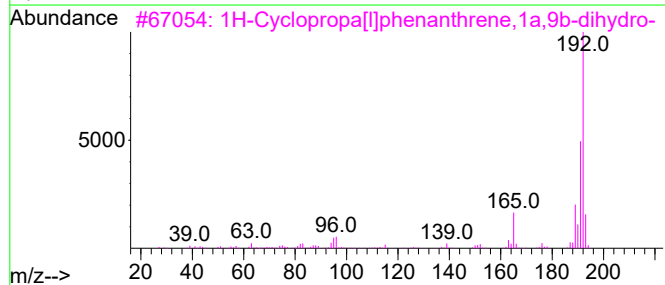
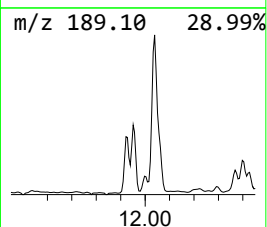
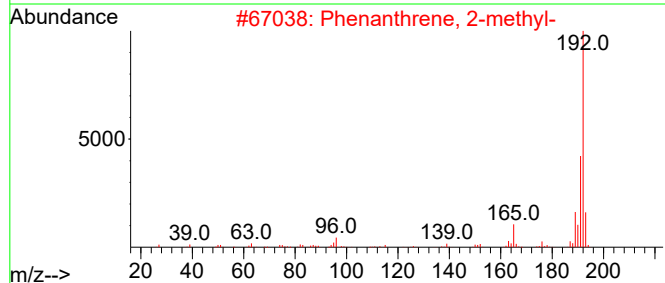
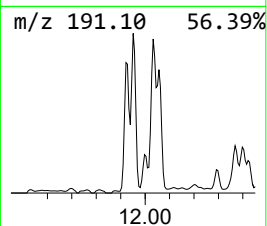
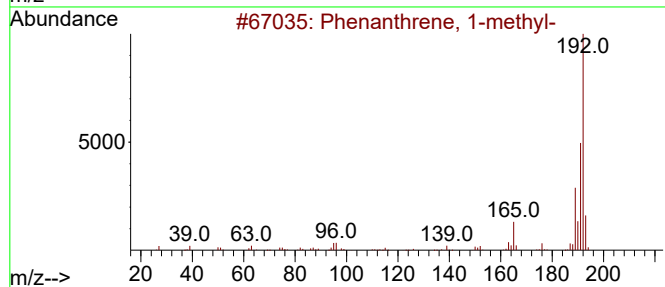
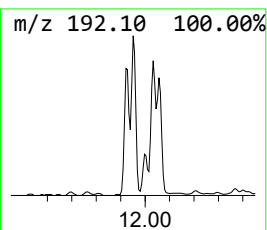
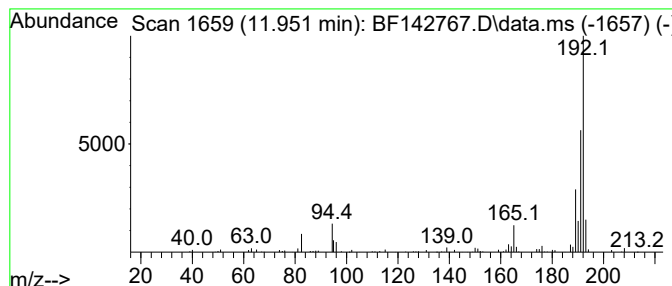
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF061125.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 5

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.951 | 4.92 ng | 103893 | Phenanthrene-d10 | 11.422 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 96   |
| 2    |      | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 96   |
| 3    |      | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 95   |
| 4    |      | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 95   |
| 5    |      | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 94   |



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Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

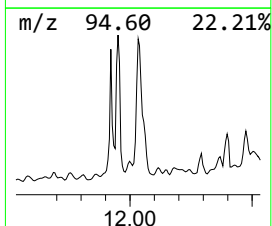
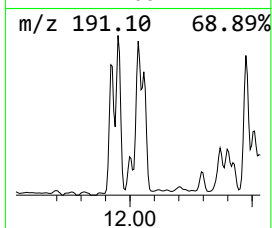
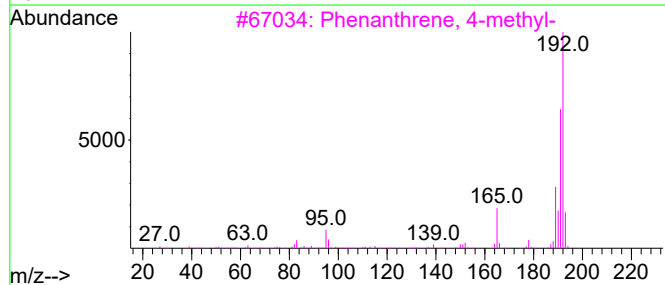
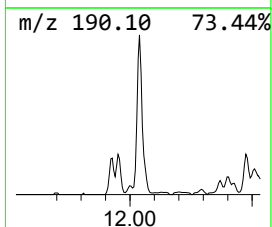
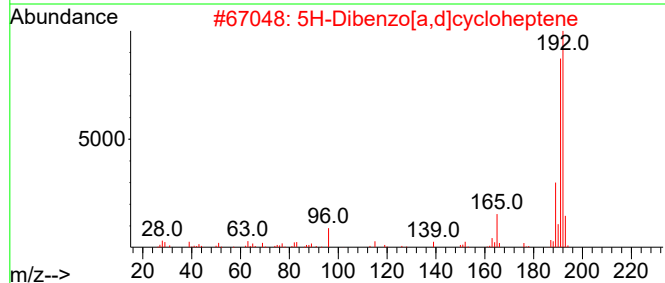
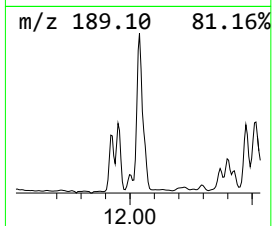
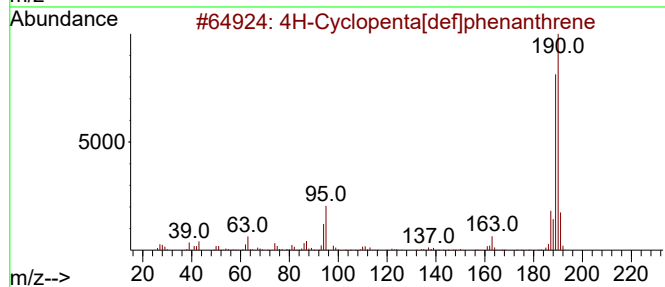
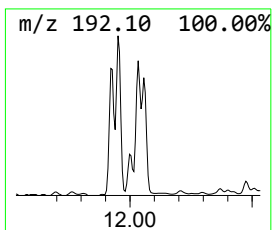
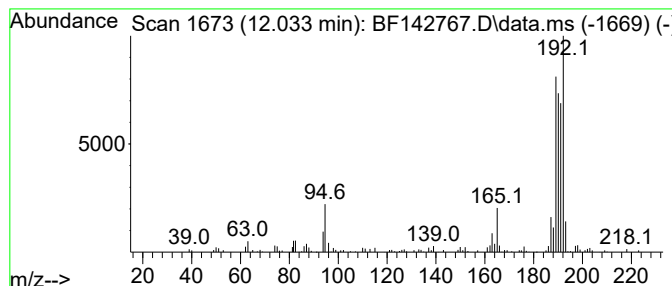
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 12.033 | 10.87 ng | 229576 | Phenanthrene-d10 | 11.422 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10  | 000203-64-5 | 50   |
| 2    |    |   | 5H-Dibenzo[a,d]cycloheptene    | 192 | C15H12  | 000256-81-5 | 46   |
| 3    |    |   | Phenanthrene, 4-methyl-        | 192 | C15H12  | 000832-64-4 | 42   |
| 4    |    |   | Anthracene, 1-methyl-          | 192 | C15H12  | 000610-48-0 | 42   |
| 5    |    |   | Phenanthrene, 1-methyl-        | 192 | C15H12  | 000832-69-9 | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
Data File : BF142767.D  
Acq On : 12 Jun 2025 19:56  
Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

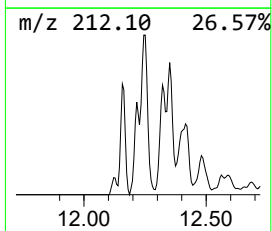
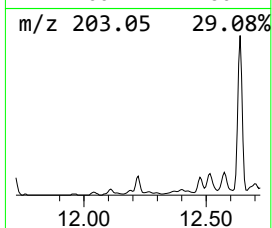
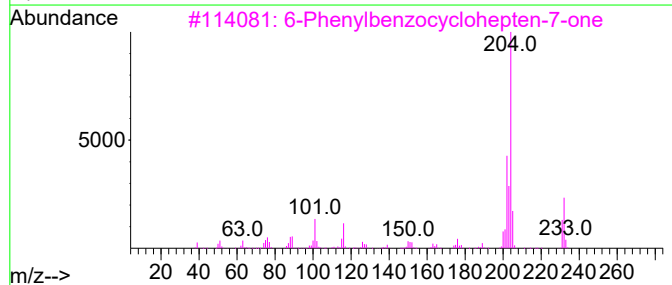
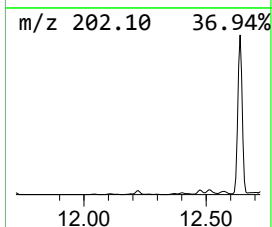
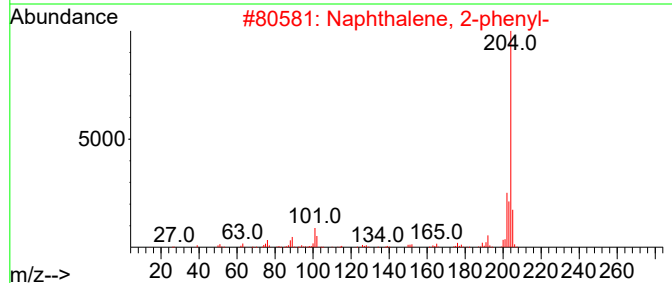
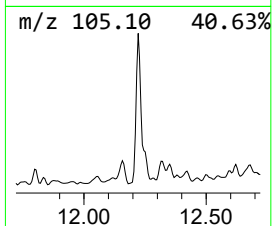
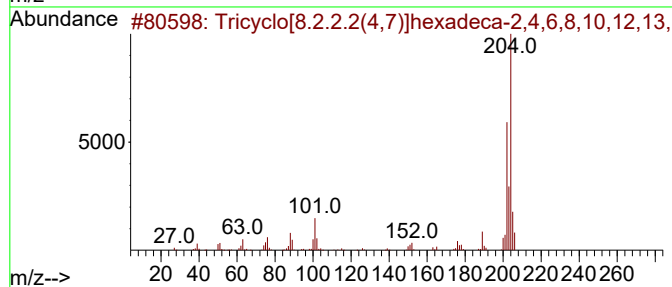
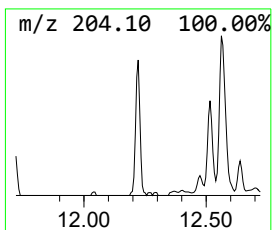
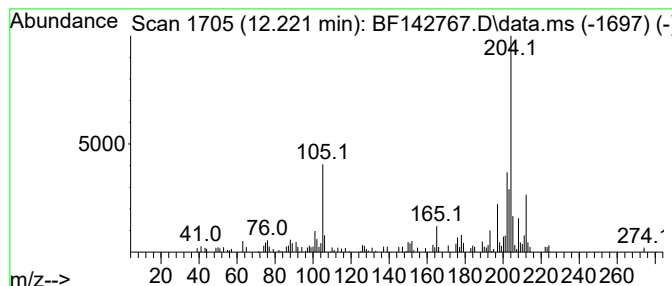
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 9

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 12.221 | 3.76 ng | 79391 | Phenanthrene-d10 | 11.422 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12  | 006572-60-7 | 70   |
| 2    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 70   |
| 3    |    |   | 6-Phenylbenzocyclohepten-7-one      | 232 | C17H12O | 093327-56-1 | 52   |
| 4    |    |   | Naphthalene, 1-phenyl-              | 204 | C16H12  | 000605-02-7 | 46   |
| 5    |    |   | 5H-Dibenzo[a,d]cycloheptene, 5-m... | 204 | C16H12  | 002975-79-3 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
Data File : BF142767.D  
Acq On : 12 Jun 2025 19:56  
Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

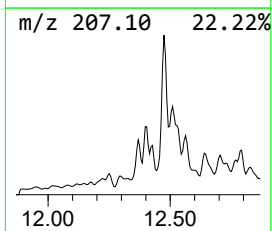
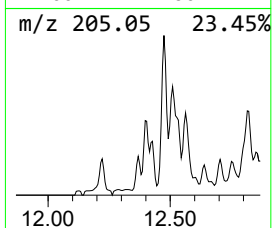
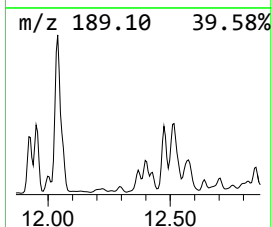
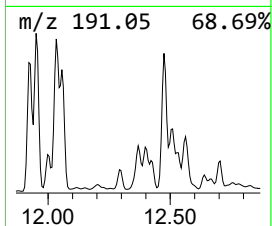
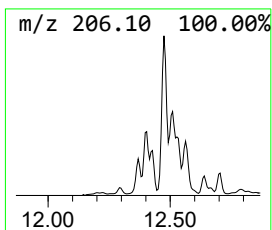
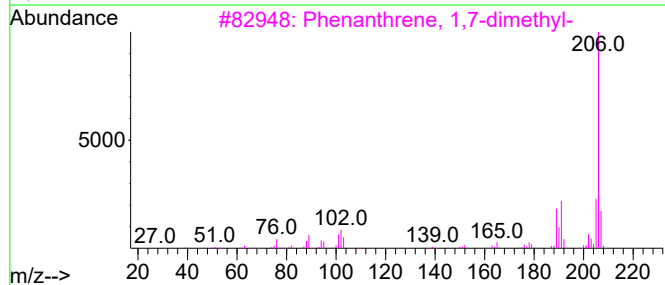
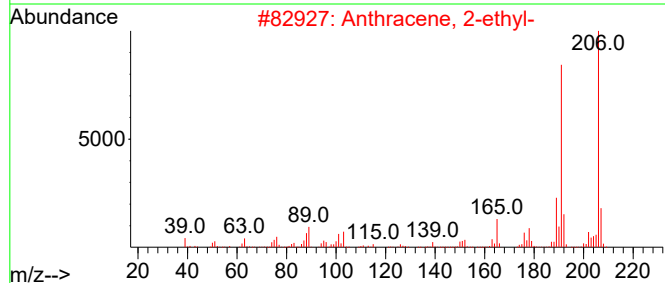
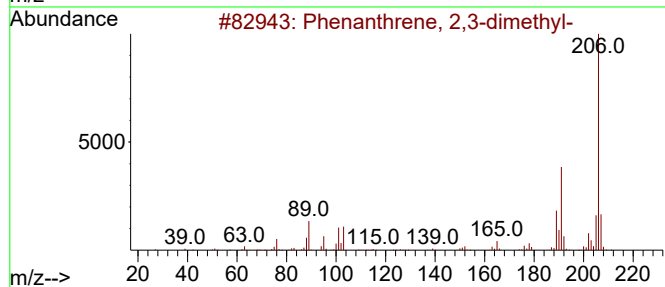
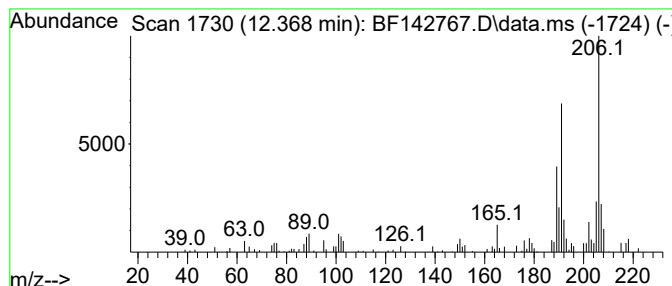
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Phenanthrene, 2,3-dimethyl- Concentration Rank 12

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 12.368 | 2.28 ng | 48130 | Phenanthrene-d10 | 11.422 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 93   |
| 2    |    |   | Anthracene, 2-ethyl-        | 206 | C16H14  | 052251-71-5 | 92   |
| 3    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 87   |
| 4    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 87   |
| 5    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
Data File : BF142767.D  
Acq On : 12 Jun 2025 19:56  
Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

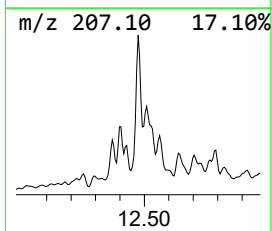
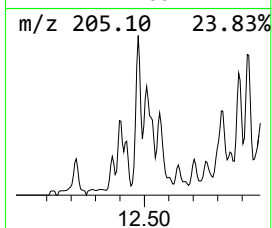
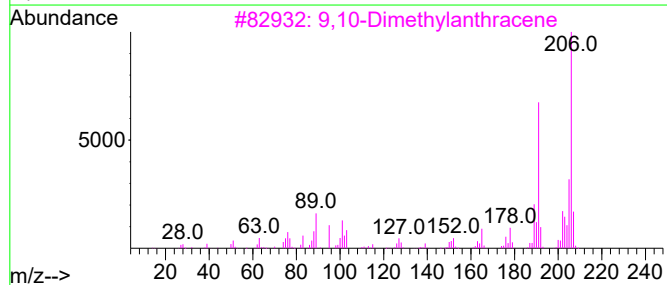
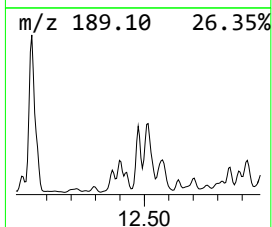
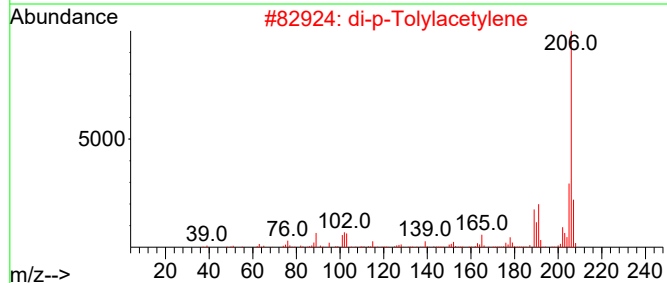
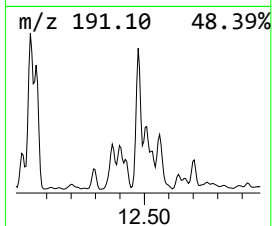
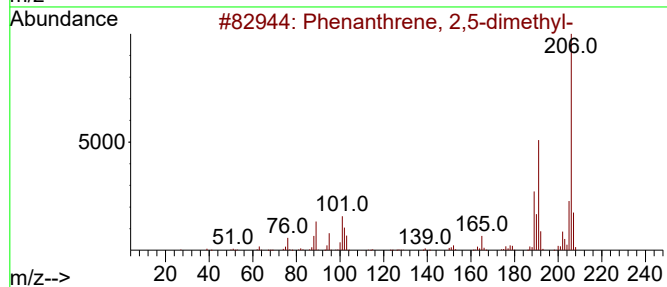
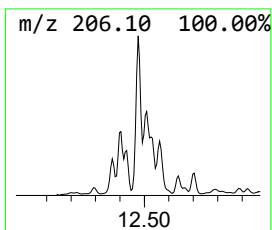
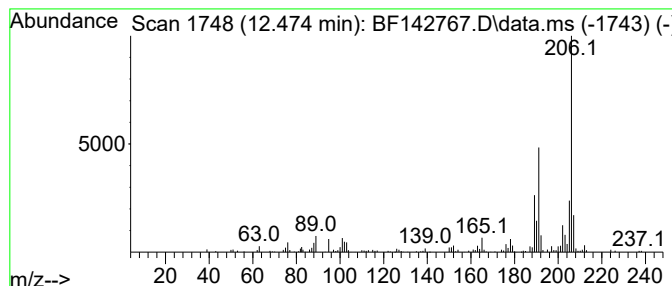
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 3

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 12.474 | 7.48 ng | 158034 | Phenanthrene-d10 | 11.422 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 96   |
| 2    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 96   |
| 3    |    |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 93   |
| 4    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 93   |
| 5    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
Data File : BF142767.D  
Acq On : 12 Jun 2025 19:56  
Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

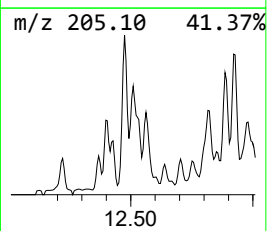
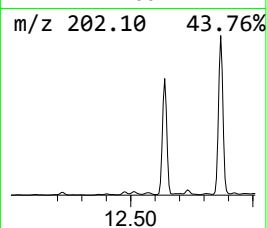
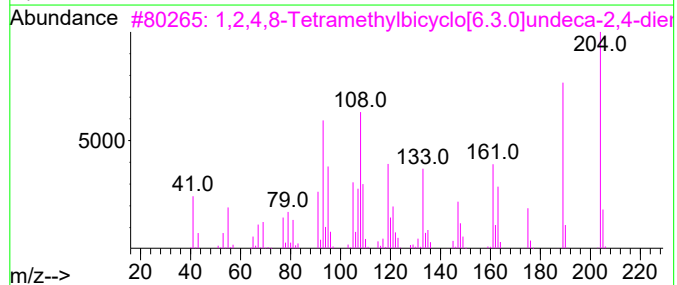
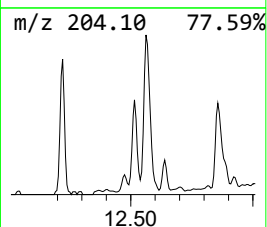
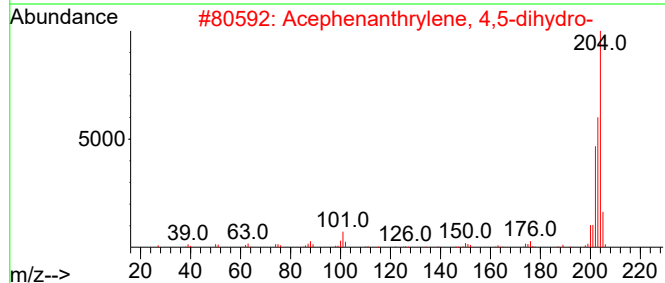
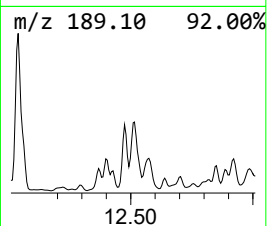
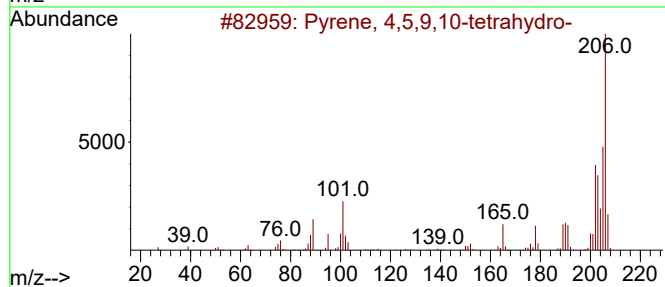
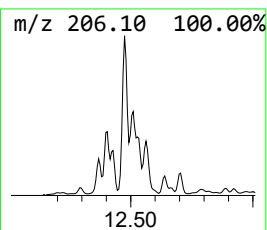
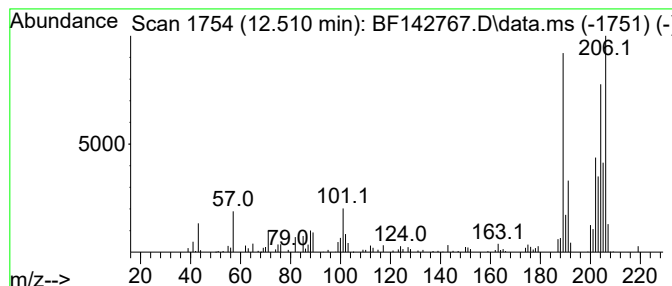
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 unknown12.510 Concentration Rank 6

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 12.510 | 4.66 ng | 98517 | Phenanthrene-d10 | 11.422 |

| Hit# | of | 5 | Tentative ID                                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 4,5,9,10-tetrahydro-                      | 206 | C16H14  | 000781-17-9 | 42   |
| 2    |    |   | Acephenanthrylene, 4,5-dihydro-                   | 204 | C16H12  | 006232-48-0 | 38   |
| 3    |    |   | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-dien- | 204 | C15H24  | 137235-51-9 | 38   |
| 4    |    |   | Naphthalene, 2-phenyl-                            | 204 | C16H12  | 000612-94-2 | 35   |
| 5    |    |   | di-p-Tolylacetylene                               | 206 | C16H14  | 002789-88-0 | 30   |



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Data File : BF142767.D  
Acq On : 12 Jun 2025 19:56  
Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

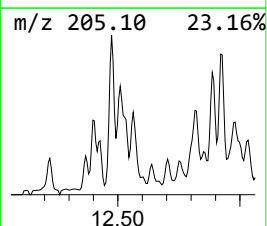
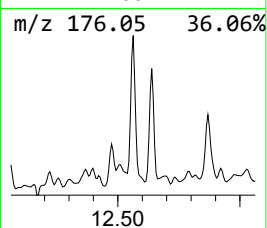
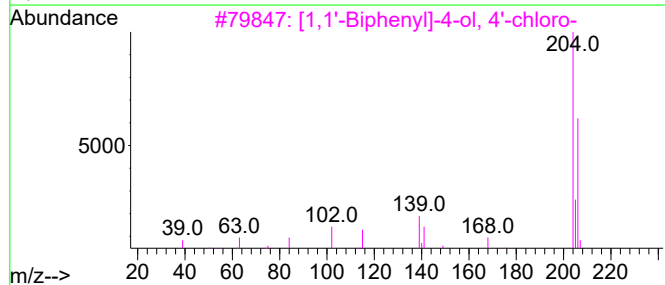
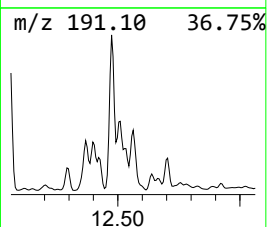
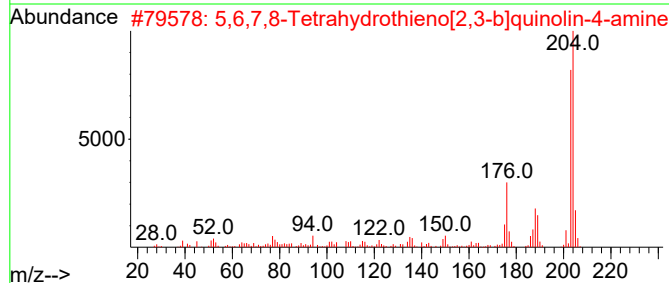
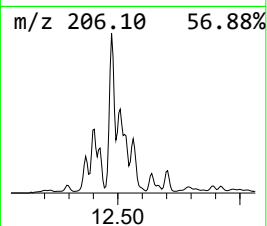
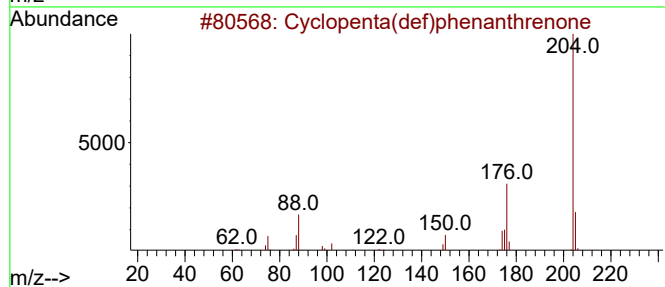
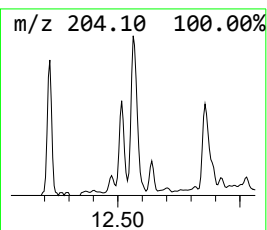
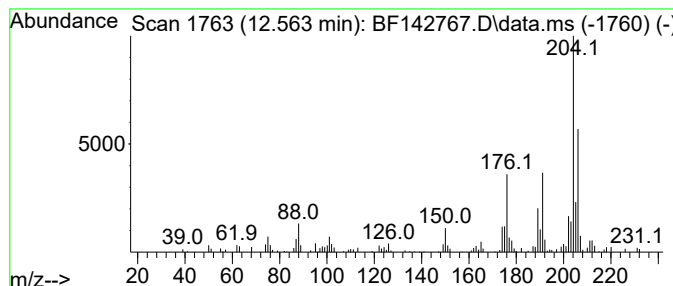
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Cyclopenta(def)phenanthrenone Concentration Rank 8

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 12.563 | 3.77 ng | 79642 | Phenanthrene-d10 | 11.422 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Cyclopenta(def)phenanthrenone       | 204 | C15H8O    | 005737-13-3 | 89   |
| 2    |    |   | 5,6,7,8-Tetrahydrothieno[2,3-b]q... | 204 | C11H12N2S | 122914-50-5 | 49   |
| 3    |    |   | [1,1'-Biphenyl]-4-ol, 4'-chloro-    | 204 | C12H9ClO  | 028034-99-3 | 49   |
| 4    |    |   | [1,1'-Biphenyl]-4,4'-dicarbonitrile | 204 | C14H8N2   | 001591-30-6 | 43   |
| 5    |    |   | 9,10-Dimethylanthracene             | 206 | C16H14    | 000781-43-1 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
Data File : BF142767.D  
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Operator : RC/JU  
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Misc :  
ALS Vial : 21 Sample Multiplier: 1

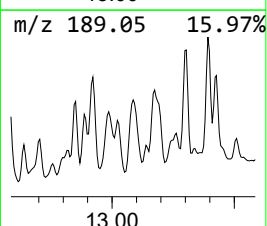
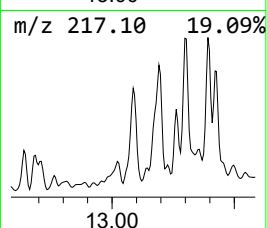
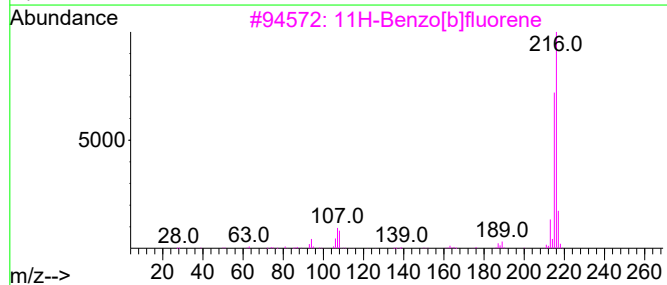
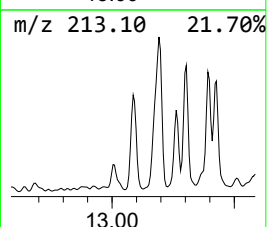
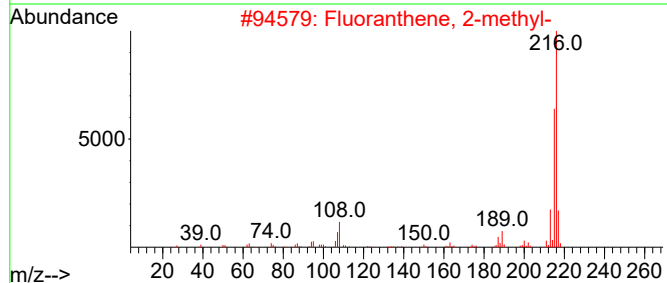
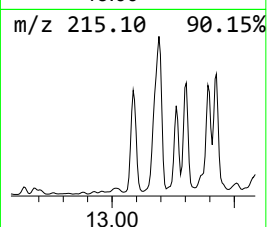
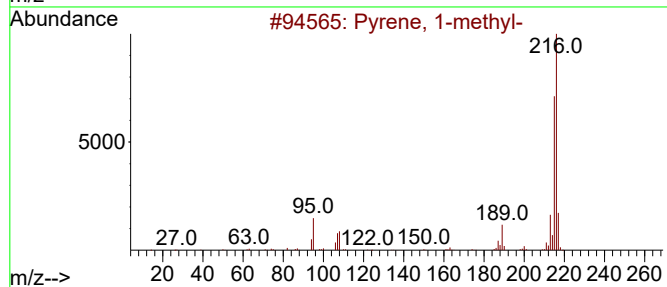
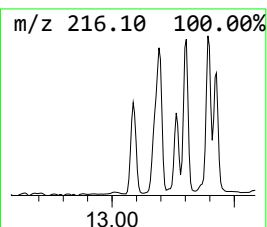
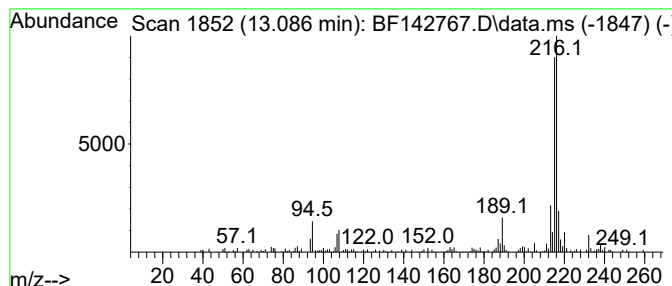
Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF061125.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 10 Pyrene, 1-methyl- Concentration Rank 10

| R.T.      | EstConc                 | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------|--------|------------------|-------------|------|
| 13.086    | 3.51 ng                 | 127356 | Chrysene-d12     | 14.068      |      |
| Hit# of 5 | Tentative ID            | MW     | MolForm          | CAS#        | Qual |
| 1         | Pyrene, 1-methyl-       | 216    | C17H12           | 002381-21-7 | 93   |
| 2         | Fluoranthene, 2-methyl- | 216    | C17H12           | 033543-31-6 | 90   |
| 3         | 11H-Benzo[b]fluorene    | 216    | C17H12           | 000243-17-4 | 87   |
| 4         | 7H-Benzo[c]fluorene     | 216    | C17H12           | 000205-12-9 | 74   |
| 5         | 11H-Benzo[a]fluorene    | 216    | C17H12           | 000238-84-6 | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
Data File : BF142767.D  
Acq On : 12 Jun 2025 19:56  
Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

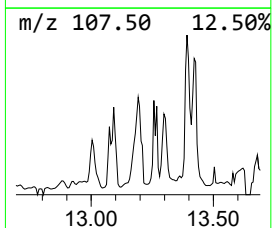
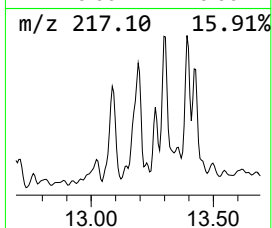
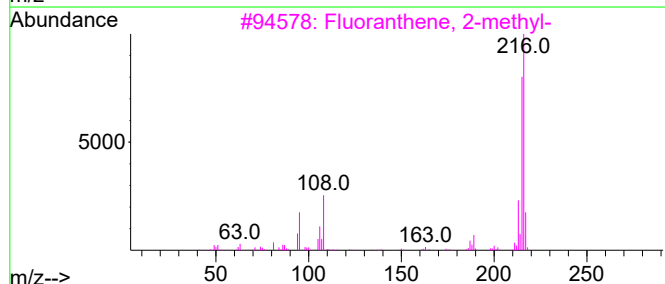
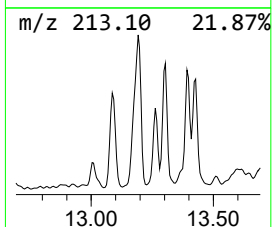
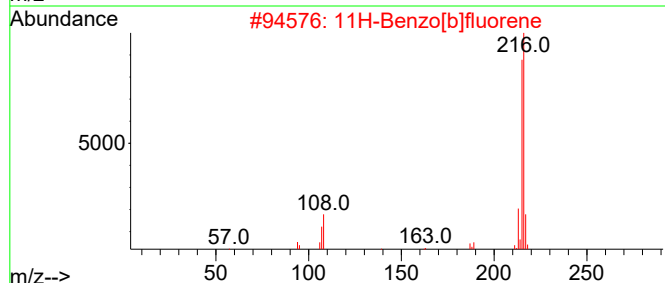
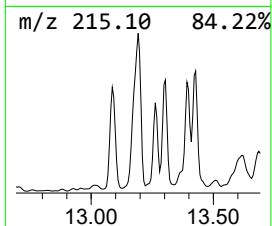
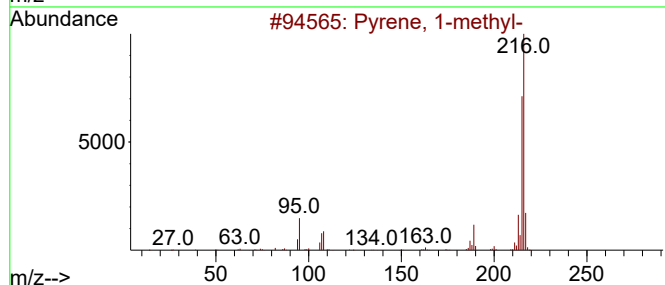
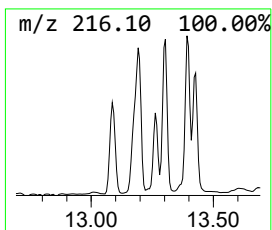
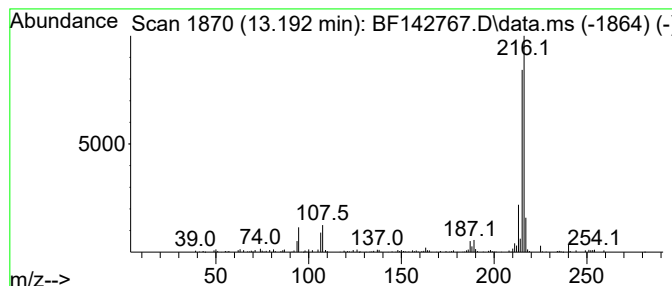
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 7

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 13.192 | 3.98 ng | 144329 | Chrysene-d12     | 14.068 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 93   |
| 2    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 91   |
| 3    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 90   |
| 4    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 81   |
| 5    |    |   | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF061225\  
Data File : BF142767.D  
Acq On : 12 Jun 2025 19:56  
Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

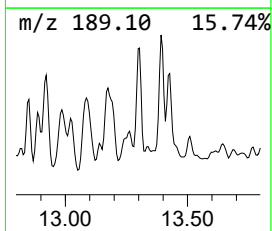
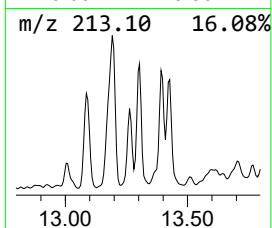
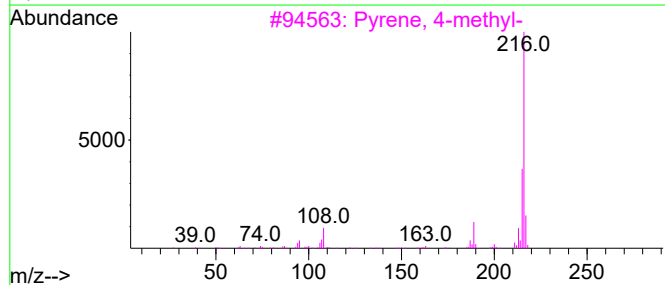
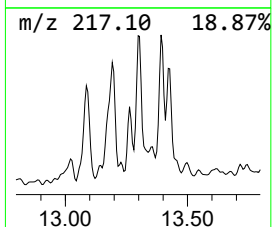
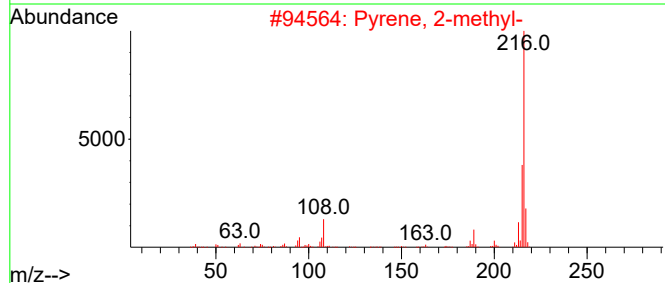
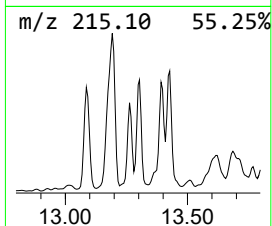
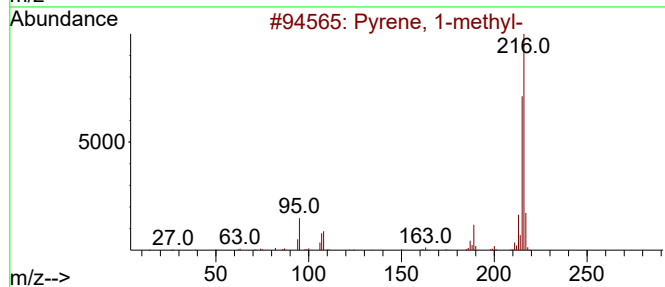
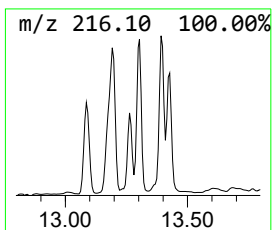
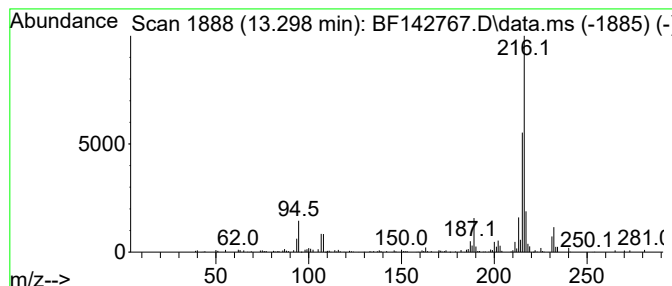
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF061125.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 Pyrene, 2-methyl- Concentration Rank 11

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 13.298 | 2.35 ng | 85163 | Chrysene-d12     | 14.068 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 95   |
| 2    |    |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 93   |
| 3    |    |   | Pyrene, 4-methyl-       | 216 | C17H12  | 003353-12-6 | 93   |
| 4    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 89   |
| 5    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF0611225\  
Data File : BF142767.D  
Acq On : 12 Jun 2025 19:56  
Operator : RC/JU  
Sample : Q2195-01 5X  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OK-01-060325

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF061125.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-... | 5.098  | 23.0    | ng    | 271146   | 1                     | 6.892  | 235974 | 20.0 |
| Anthracene, 2-m... | 11.927 | 4.9     | ng    | 104404   | 4                     | 11.422 | 422475 | 20.0 |
| Phenanthrene, 1... | 11.951 | 4.9     | ng    | 103893   | 4                     | 11.422 | 422475 | 20.0 |
| 4H-Cyclopenta[d... | 12.033 | 10.9    | ng    | 229576   | 4                     | 11.422 | 422475 | 20.0 |
| Tricyclo[8.2.2...  | 12.221 | 3.8     | ng    | 79391    | 4                     | 11.422 | 422475 | 20.0 |
| Phenanthrene, 2... | 12.368 | 2.3     | ng    | 48130    | 4                     | 11.422 | 422475 | 20.0 |
| Phenanthrene, 2... | 12.474 | 7.5     | ng    | 158034   | 4                     | 11.422 | 422475 | 20.0 |
| unknown12.510      | 12.510 | 4.7     | ng    | 98517    | 4                     | 11.422 | 422475 | 20.0 |
| Cyclopenta(def)... | 12.563 | 3.8     | ng    | 79642    | 4                     | 11.422 | 422475 | 20.0 |
| Pyrene, 1-methyl-  | 13.086 | 3.5     | ng    | 127356   | 5                     | 14.068 | 725475 | 20.0 |
| 11H-Benzo[b]flu... | 13.192 | 4.0     | ng    | 144329   | 5                     | 14.068 | 725475 | 20.0 |
| Pyrene, 2-methyl-  | 13.298 | 2.4     | ng    | 85163    | 5                     | 14.068 | 725475 | 20.0 |