

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
 Data File : BF139329.D  
 Acq On : 11 Sep 2024 16:50  
 Operator : RC/JU  
 Sample : P3912-01  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 282-COMPOSITE-1

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-BF090424.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139329.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         | 2.181       | 7             | 15          | 21           | rVB      | 2628476        | 4231509       | 98.42%          | 9.023%        |
| 2         | 2.999       | 143           | 154         | 163          | rBV      | 125227         | 226199        | 5.26%           | 0.482%        |
| 3         | 4.475       | 399           | 405         | 412          | rBV      | 113132         | 158543        | 3.69%           | 0.338%        |
| 4         | 5.093       | 502           | 510         | 516          | rVB      | 58581          | 89579         | 2.08%           | 0.191%        |
| 5         | 5.498       | 573           | 579         | 584          | rBV      | 1034552        | 1387737       | 32.28%          | 2.959%        |
| 6         | 6.504       | 744           | 750         | 757          | rVB      | 858627         | 1140738       | 26.53%          | 2.432%        |
| 7         | 6.704       | 777           | 784         | 789          | rBV      | 36783          | 52911         | 1.23%           | 0.113%        |
| 8         | 6.881       | 809           | 814         | 819          | rBV      | 528455         | 640020        | 14.89%          | 1.365%        |
| 9         | 7.440       | 901           | 909         | 914          | rBV      | 819179         | 1054085       | 24.52%          | 2.248%        |
| 10        | 7.692       | 947           | 952         | 959          | rVB2     | 45459          | 61360         | 1.43%           | 0.131%        |
| 11        | 7.898       | 980           | 987         | 992          | rVB      | 51575          | 75141         | 1.75%           | 0.160%        |
| 12        | 8.187       | 1033          | 1036        | 1041         | rVB      | 1040259        | 1343528       | 31.25%          | 2.865%        |
| 13        | 8.234       | 1041          | 1044        | 1049         | rVB      | 79782          | 99602         | 2.32%           | 0.212%        |
| 14        | 8.475       | 1081          | 1085        | 1094         | rVV      | 213292         | 332513        | 7.73%           | 0.709%        |
| 15        | 8.545       | 1094          | 1097        | 1103         | rVB3     | 42530          | 60518         | 1.41%           | 0.129%        |
| 16        | 8.751       | 1128          | 1132        | 1136         | rVB      | 89755          | 105273        | 2.45%           | 0.224%        |
| 17        | 8.828       | 1141          | 1145        | 1148         | rVB2     | 38268          | 47335         | 1.10%           | 0.101%        |
| 18        | 8.881       | 1148          | 1154        | 1158         | rBV      | 1489575        | 1944237       | 45.22%          | 4.146%        |
| 19        | 8.928       | 1158          | 1162        | 1165         | rBV2     | 36868          | 47461         | 1.10%           | 0.101%        |
| 20        | 8.975       | 1165          | 1170        | 1175         | rVB      | 1022953        | 1278620       | 29.74%          | 2.726%        |
| 21        | 9.239       | 1209          | 1215        | 1220         | rVB      | 1489698        | 1883131       | 43.80%          | 4.015%        |
| 22        | 9.339       | 1225          | 1232        | 1237         | rVV      | 332778         | 419173        | 9.75%           | 0.894%        |
| 23        | 9.434       | 1241          | 1248        | 1254         | rVB      | 132988         | 202338        | 4.71%           | 0.431%        |
| 24        | 9.498       | 1254          | 1259        | 1265         | rBV      | 295225         | 464943        | 10.81%          | 0.991%        |
| 25        | 9.575       | 1267          | 1272        | 1275         | rVV      | 497650         | 711783        | 16.56%          | 1.518%        |
| 26        | 9.604       | 1275          | 1277        | 1281         | rVB      | 275972         | 279662        | 6.50%           | 0.596%        |
| 27        | 9.692       | 1288          | 1292        | 1298         | rBV      | 168989         | 275285        | 6.40%           | 0.587%        |
| 28        | 9.775       | 1302          | 1306        | 1311         | rVB      | 166934         | 216330        | 5.03%           | 0.461%        |
| 29        | 9.916       | 1323          | 1330        | 1333         | rBV2     | 710564         | 1178061       | 27.40%          | 2.512%        |
| 30        | 9.957       | 1333          | 1337        | 1341         | rVV      | 2816289        | 3721489       | 86.56%          | 7.935%        |
| 31        | 10.016      | 1345          | 1347        | 1351         | rVV3     | 47391          | 72548         | 1.69%           | 0.155%        |
| 32        | 10.075      | 1351          | 1357        | 1361         | rVV2     | 142000         | 234687        | 5.46%           | 0.500%        |
| 33        | 10.122      | 1361          | 1365        | 1370         | rVV      | 1580582        | 2090589       | 48.62%          | 4.458%        |
| 34        | 10.157      | 1370          | 1371        | 1374         | rVV      | 51431          | 47725         | 1.11%           | 0.102%        |
| 35        | 10.192      | 1374          | 1377        | 1380         | rVV      | 65672          | 78128         | 1.82%           | 0.167%        |
| 36        | 10.234      | 1380          | 1384        | 1387         | rVV2     | 60077          | 78347         | 1.82%           | 0.167%        |

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 Sample : P3912-01  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 282-COMPOSITE-1

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-BF090424.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 37 | 10.322 | 1394 | 1399 | 1401 | rBV  | 63615   | 91541   | 2.13%   | 0.195% |
| 38 | 10.345 | 1401 | 1403 | 1409 | rVB2 | 64502   | 70462   | 1.64%   | 0.150% |
| 39 | 10.416 | 1409 | 1415 | 1418 | rBV2 | 44669   | 72913   | 1.70%   | 0.155% |
| 40 | 10.469 | 1418 | 1424 | 1429 | rBV  | 2535177 | 3312124 | 77.04%  | 7.062% |
| 41 | 10.551 | 1436 | 1438 | 1444 | rVB2 | 144218  | 155082  | 3.61%   | 0.331% |
| 42 | 10.622 | 1444 | 1450 | 1454 | rVB2 | 257016  | 440013  | 10.23%  | 0.938% |
| 43 | 10.704 | 1454 | 1464 | 1469 | rBV  | 1382538 | 1911107 | 44.45%  | 4.075% |
| 44 | 10.751 | 1469 | 1472 | 1476 | rVB  | 57442   | 74938   | 1.74%   | 0.160% |
| 45 | 10.828 | 1478 | 1485 | 1492 | rBV  | 187078  | 295206  | 6.87%   | 0.629% |
| 46 | 10.892 | 1492 | 1496 | 1500 | rBV4 | 50131   | 81175   | 1.89%   | 0.173% |
| 47 | 11.010 | 1509 | 1516 | 1519 | rBV2 | 226717  | 380828  | 8.86%   | 0.812% |
| 48 | 11.039 | 1519 | 1521 | 1524 | rVV  | 82826   | 101561  | 2.36%   | 0.217% |
| 49 | 11.092 | 1527 | 1530 | 1534 | rVV  | 87297   | 113880  | 2.65%   | 0.243% |
| 50 | 11.163 | 1534 | 1542 | 1546 | rVV4 | 54881   | 146238  | 3.40%   | 0.312% |
| 51 | 11.210 | 1546 | 1550 | 1554 | rVB  | 101219  | 129426  | 3.01%   | 0.276% |
| 52 | 11.257 | 1554 | 1558 | 1561 | rBV2 | 60878   | 112349  | 2.61%   | 0.240% |
| 53 | 11.298 | 1561 | 1565 | 1570 | rVB  | 330051  | 431805  | 10.04%  | 0.921% |
| 54 | 11.433 | 1584 | 1588 | 1593 | rVB  | 3055506 | 4299456 | 100.00% | 9.168% |
| 55 | 11.480 | 1593 | 1596 | 1599 | rVV  | 250014  | 280905  | 6.53%   | 0.599% |
| 56 | 11.510 | 1599 | 1601 | 1606 | rVB2 | 51880   | 61093   | 1.42%   | 0.130% |
| 57 | 11.633 | 1618 | 1622 | 1628 | rBV  | 129805  | 186856  | 4.35%   | 0.398% |
| 58 | 11.698 | 1629 | 1633 | 1636 | rBV2 | 53033   | 71912   | 1.67%   | 0.153% |
| 59 | 11.792 | 1646 | 1649 | 1657 | rVB4 | 40146   | 81565   | 1.90%   | 0.174% |
| 60 | 11.904 | 1664 | 1668 | 1669 | rBV  | 177716  | 188825  | 4.39%   | 0.403% |
| 61 | 11.933 | 1669 | 1673 | 1678 | rVV  | 856546  | 1151369 | 26.78%  | 2.455% |
| 62 | 11.980 | 1678 | 1681 | 1683 | rVV  | 80057   | 106289  | 2.47%   | 0.227% |
| 63 | 12.022 | 1683 | 1688 | 1695 | rVB  | 461165  | 707077  | 16.45%  | 1.508% |
| 64 | 12.198 | 1715 | 1718 | 1720 | rBV  | 72007   | 92212   | 2.14%   | 0.197% |
| 65 | 12.539 | 1773 | 1776 | 1781 | rVB  | 34578   | 46908   | 1.09%   | 0.100% |
| 66 | 12.616 | 1782 | 1789 | 1795 | rBV  | 814028  | 1080529 | 25.13%  | 2.304% |
| 67 | 12.716 | 1801 | 1806 | 1813 | rBV2 | 161846  | 241076  | 5.61%   | 0.514% |
| 68 | 12.845 | 1821 | 1828 | 1832 | rBV  | 404499  | 641253  | 14.91%  | 1.367% |
| 69 | 12.886 | 1832 | 1835 | 1842 | rVB3 | 62995   | 93880   | 2.18%   | 0.200% |
| 70 | 12.986 | 1847 | 1852 | 1858 | rVB  | 1459694 | 1916872 | 44.58%  | 4.087% |
| 71 | 13.139 | 1874 | 1878 | 1881 | rBV  | 38898   | 55098   | 1.28%   | 0.117% |
| 72 | 13.169 | 1881 | 1883 | 1888 | rVB  | 66105   | 77622   | 1.81%   | 0.166% |
| 73 | 13.239 | 1891 | 1895 | 1898 | rBV  | 58498   | 72447   | 1.69%   | 0.154% |
| 74 | 13.816 | 1986 | 1993 | 2000 | rVB5 | 25067   | 55433   | 1.29%   | 0.118% |
| 75 | 13.904 | 2004 | 2008 | 2020 | rVB7 | 24785   | 65778   | 1.53%   | 0.140% |
| 76 | 14.039 | 2025 | 2031 | 2044 | rVV  | 451171  | 655417  | 15.24%  | 1.398% |

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Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

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-BF090424.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |       |        |
|----|--------|------|------|------|------|--------|--------|-------|--------|
| 77 | 15.516 | 2275 | 2282 | 2295 | rVB2 | 252140 | 416531 | 9.69% | 0.888% |
|----|--------|------|------|------|------|--------|--------|-------|--------|

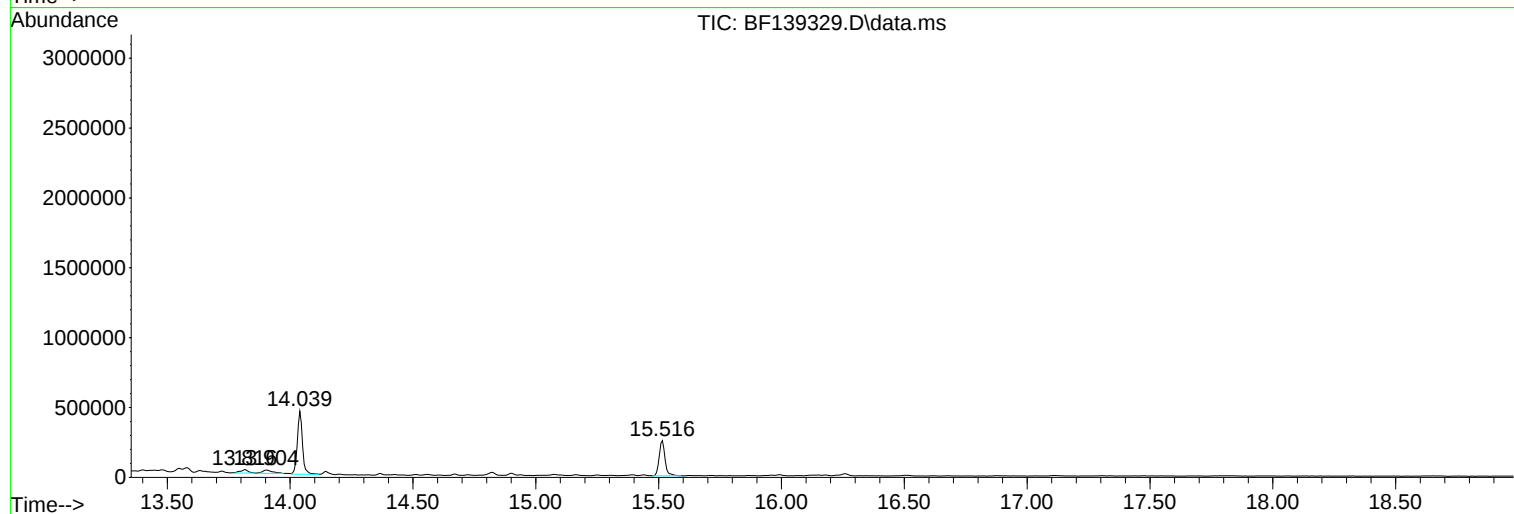
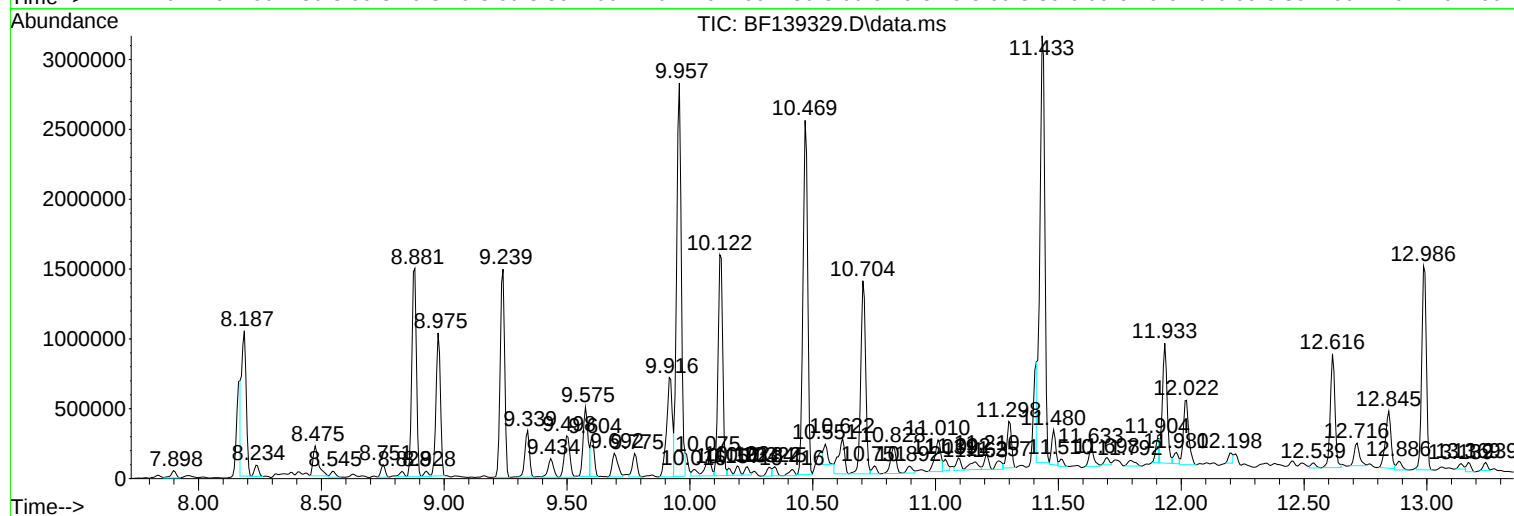
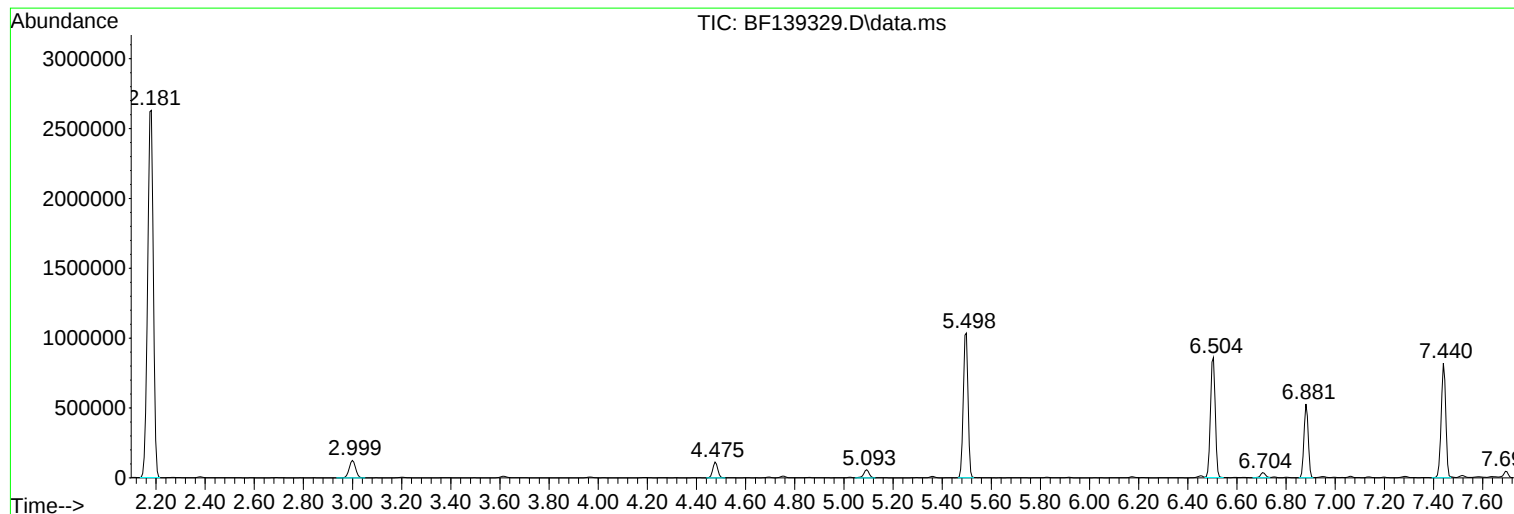
Sum of corrected areas: 46898179

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
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Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.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\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

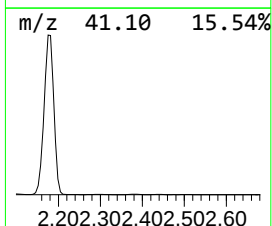
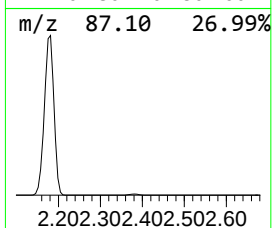
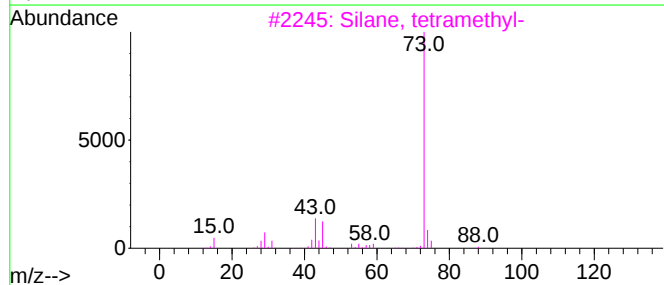
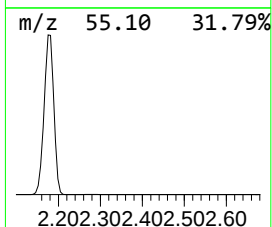
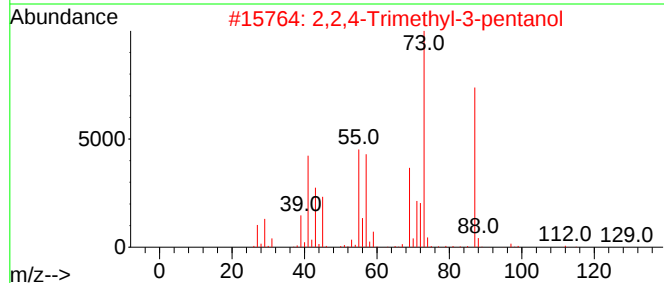
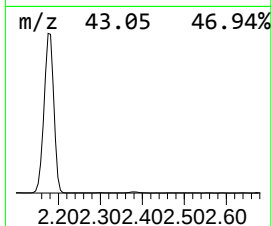
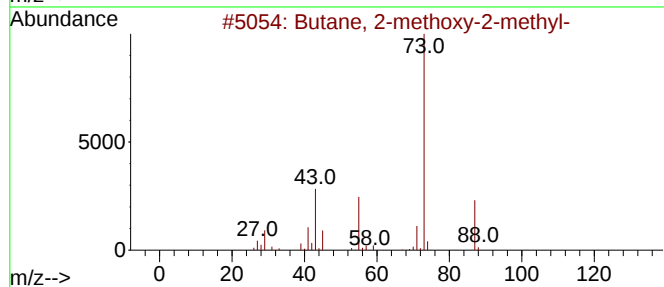
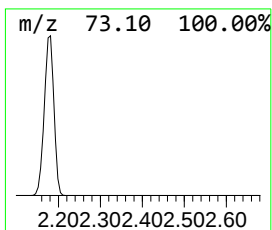
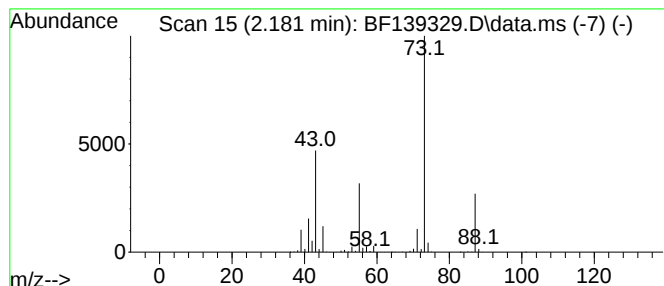
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

| R.T.  | EstConc   | Area    | Relative to ISTD       | R.T.  |
|-------|-----------|---------|------------------------|-------|
| 2.181 | 132.23 ng | 4231510 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 38   |
| 4    |    |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 23   |
| 5    |    |   | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 9    |



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ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

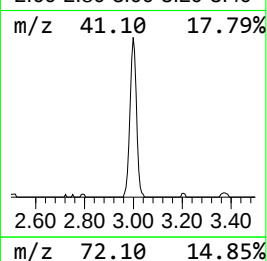
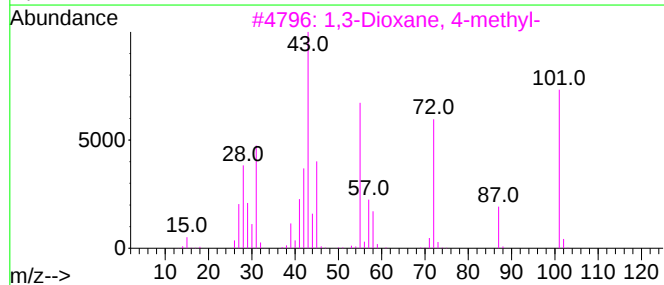
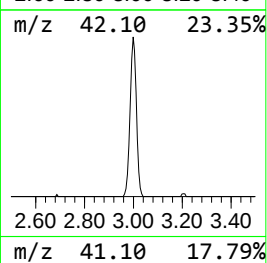
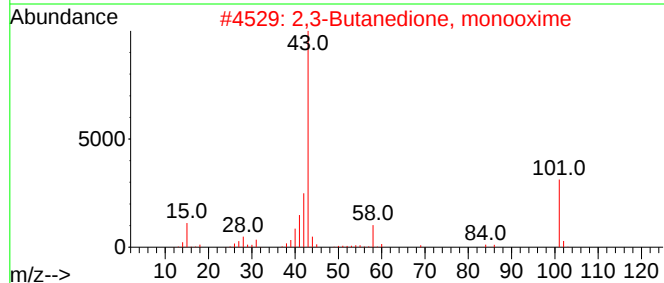
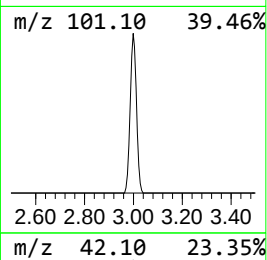
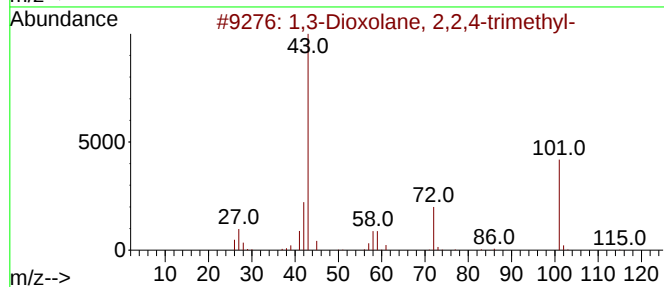
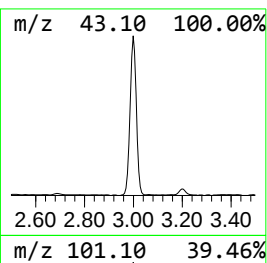
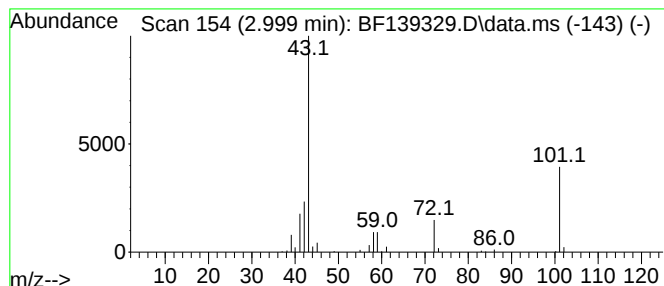
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 1,3-Dioxolane, 2,2,4-trimet... Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 2.999 | 7.07 ng | 226199 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1,3-Dioxolane, 2,2,4-trimethyl-     | 116 | C6H12O2 | 001193-11-9 | 78   |
| 2    |    |   | 2,3-Butanedione, monooxime          | 101 | C4H7NO2 | 000057-71-6 | 59   |
| 3    |    |   | 1,3-Dioxane, 4-methyl-              | 102 | C5H10O2 | 001120-97-4 | 42   |
| 4    |    |   | 1,3-Dioxolane-2-methanol, 2,4-di... | 132 | C6H12O3 | 053951-43-2 | 33   |
| 5    |    |   | Isopropyl isothiocyanate            | 101 | C4H7NS  | 002253-73-8 | 9    |



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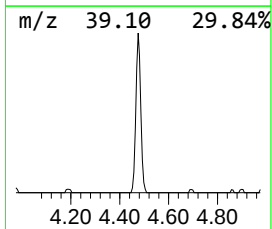
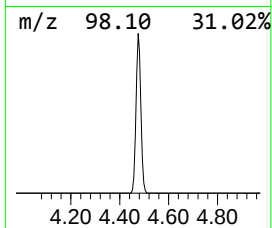
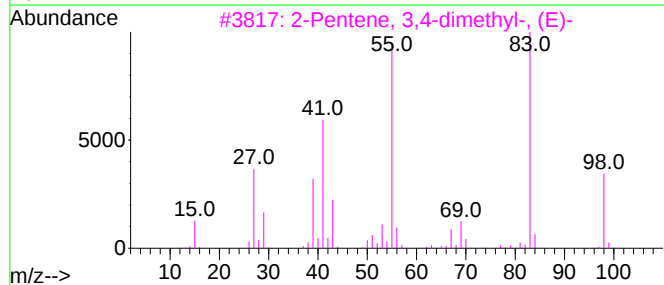
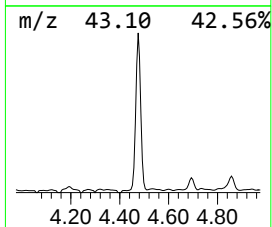
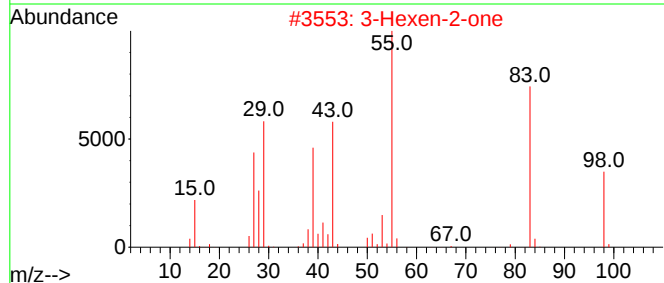
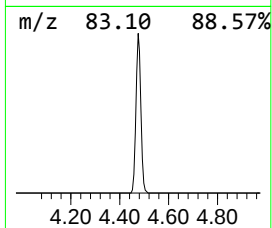
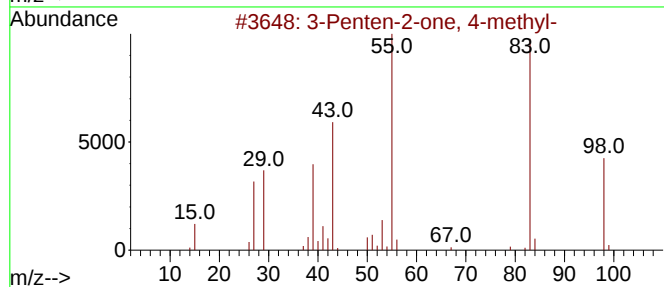
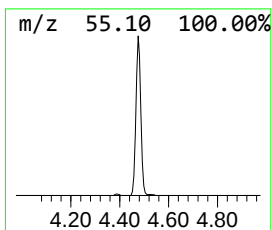
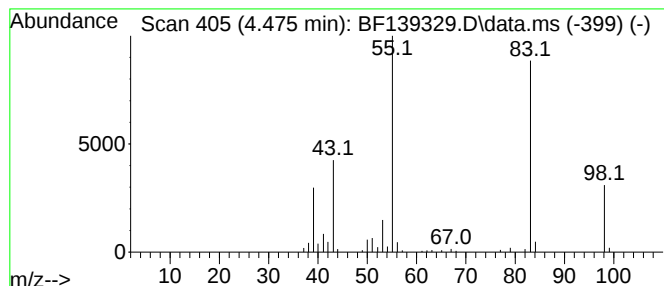
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ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 3-Penten-2-one, 4-methyl- Concentration Rank 7

| R.T.      | EstConc                        | Area   | Relative to ISTD       |         |             | R.T.  |
|-----------|--------------------------------|--------|------------------------|---------|-------------|-------|
| 4.475     | 4.95 ng                        | 158543 | 1,4-Dichlorobenzene-d4 |         |             | 6.881 |
| Hit# of 5 | Tentative ID                   |        | MW                     | MolForm | CAS#        | Qual  |
| 1         | 3-Penten-2-one, 4-methyl-      |        | 98                     | C6H10O  | 000141-79-7 | 91    |
| 2         | 3-Hexen-2-one                  |        | 98                     | C6H10O  | 000763-93-9 | 91    |
| 3         | 2-Pentene, 3,4-dimethyl-, (E)- |        | 98                     | C7H14   | 004914-92-5 | 90    |
| 4         | 2-Pentene, 3,4-dimethyl-, (Z)- |        | 98                     | C7H14   | 004914-91-4 | 86    |
| 5         | 2-Pentene, 3,4-dimethyl-       |        | 98                     | C7H14   | 024910-63-2 | 86    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

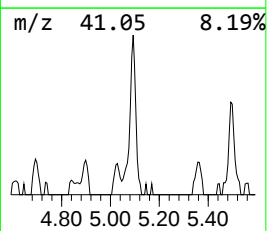
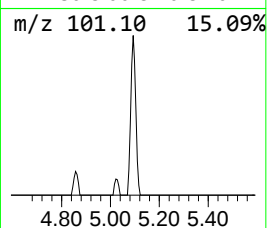
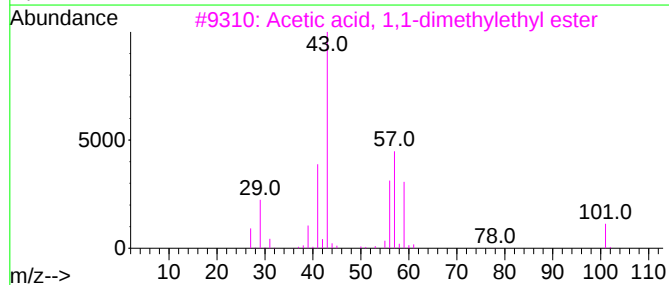
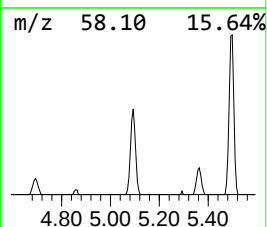
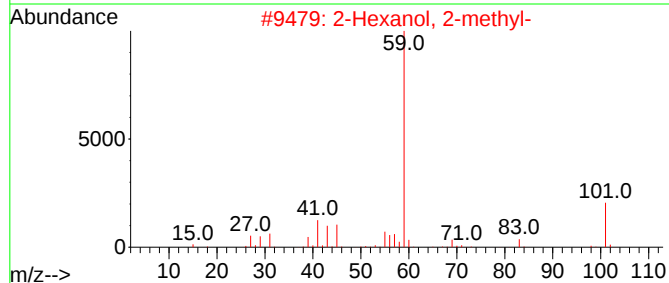
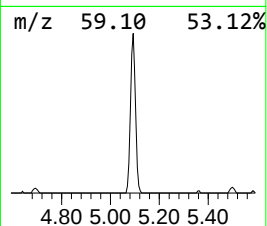
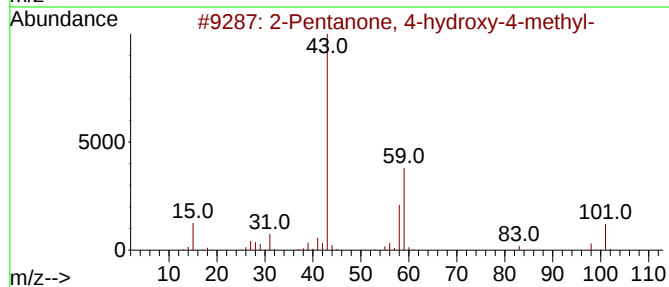
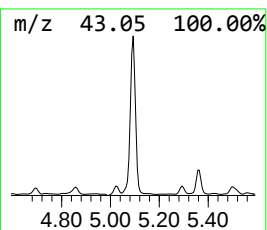
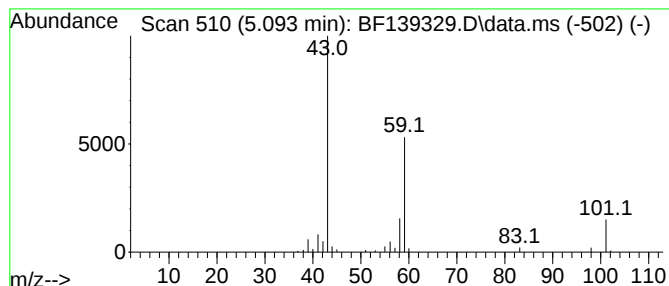
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD       | R.T.  |
|-------|---------|-------|------------------------|-------|
| 5.093 | 2.80 ng | 89579 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2 | 59   |
| 2    |      | 2-Hexanol, 2-methyl-                | 116 | C7H16O  | 000625-23-0 | 33   |
| 3    |      | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 28   |
| 4    |      | 1-Propen-2-ol, acetate              | 100 | C5H8O2  | 000108-22-5 | 12   |
| 5    |      | 2,3-Butanedione, monooxime          | 101 | C4H7NO2 | 000057-71-6 | 10   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

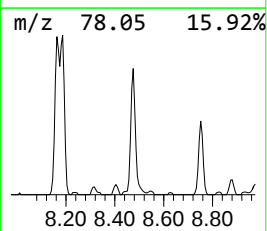
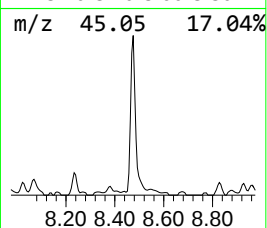
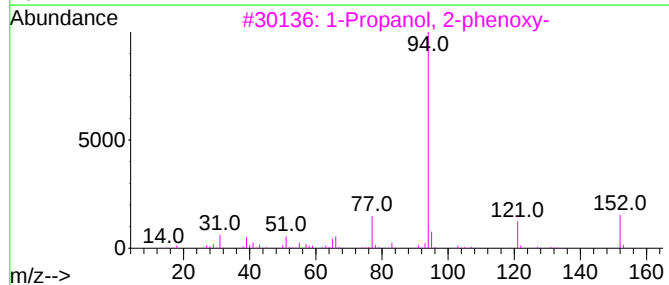
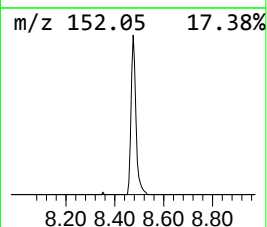
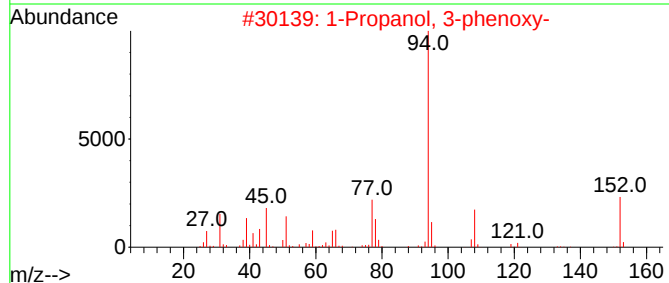
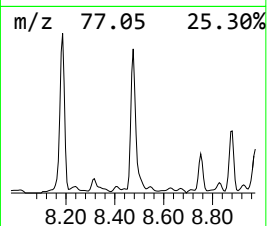
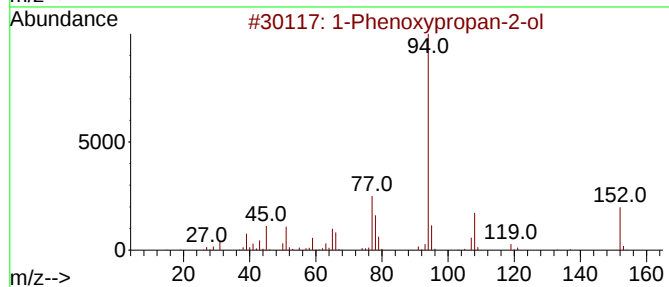
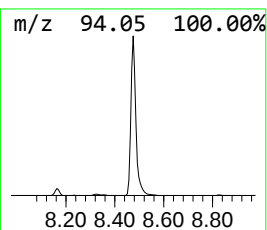
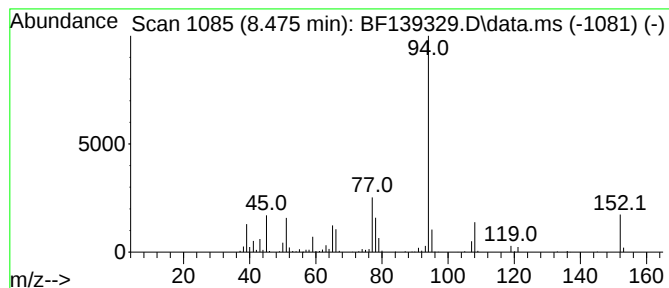
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 1-Phenoxypropan-2-ol Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 8.475 | 4.95 ng | 332513 | Naphthalene-d8   | 8.163 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Phenoxypropan-2-ol             | 152 | C9H12O2 | 000770-35-4 | 97   |
| 2    |    |   | 1-Propanol, 3-phenoxy-           | 152 | C9H12O2 | 006180-61-6 | 87   |
| 3    |    |   | 1-Propanol, 2-phenoxy-           | 152 | C9H12O2 | 004169-04-4 | 64   |
| 4    |    |   | Ethanol, 2-phenoxy-              | 138 | C8H10O2 | 000122-99-6 | 64   |
| 5    |    |   | 3-(.alpha.-Hydroxyethyl)-aniline | 137 | C8H11NO | 002454-37-7 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

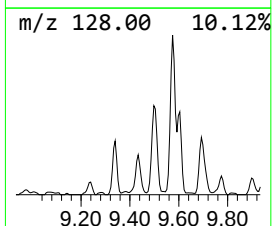
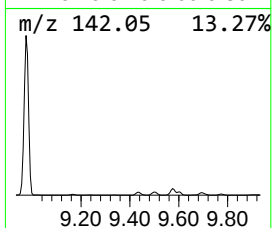
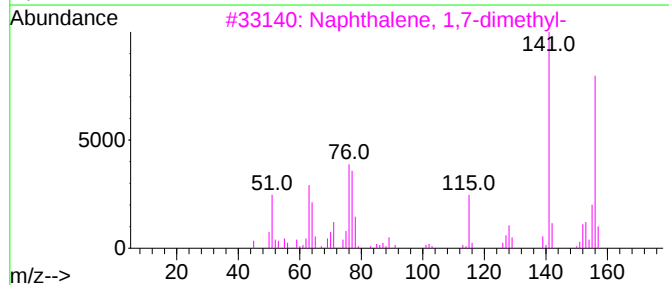
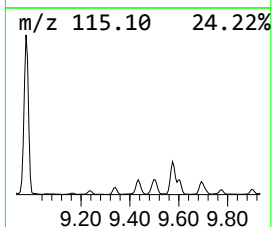
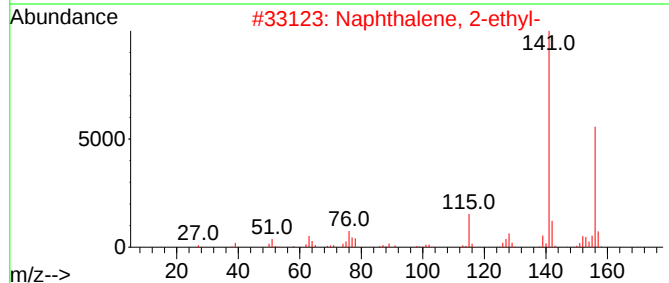
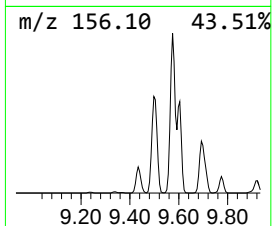
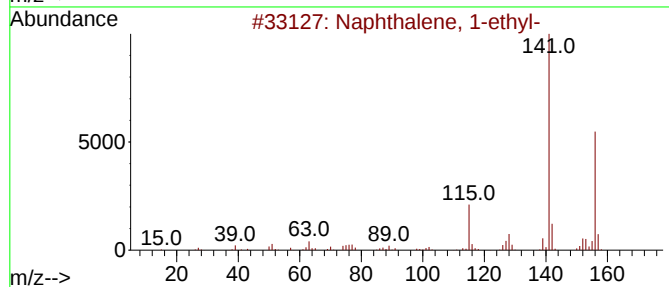
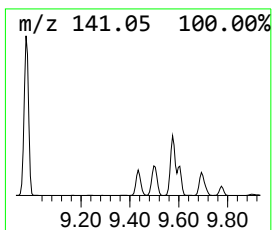
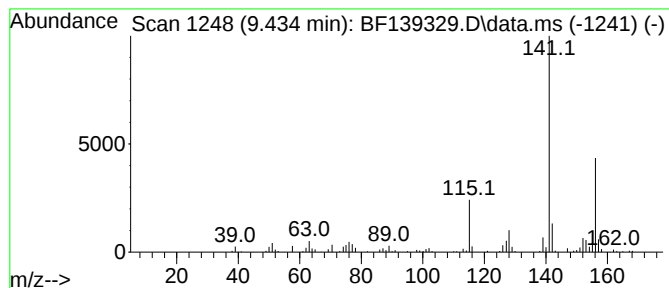
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Naphthalene, 1-ethyl- Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.434 | 3.44 ng | 202338 | Acenaphthene-d10 | 9.922 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1-ethyl-               | 156 | C12H12  | 001127-76-0 | 96   |
| 2    |    |   | Naphthalene, 2-ethyl-               | 156 | C12H12  | 000939-27-5 | 95   |
| 3    |    |   | Naphthalene, 1,7-dimethyl-          | 156 | C12H12  | 000575-37-1 | 86   |
| 4    |    |   | 2-Thiophenecarboxylic acid, 5-et... | 156 | C7H8O2S | 023229-72-3 | 64   |
| 5    |    |   | 2,4-Difluoromesitylene              | 156 | C9H10F2 | 000392-61-0 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

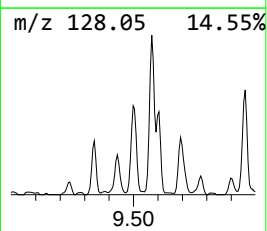
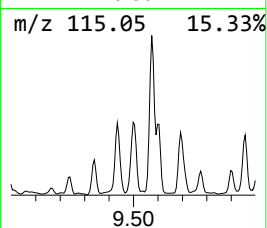
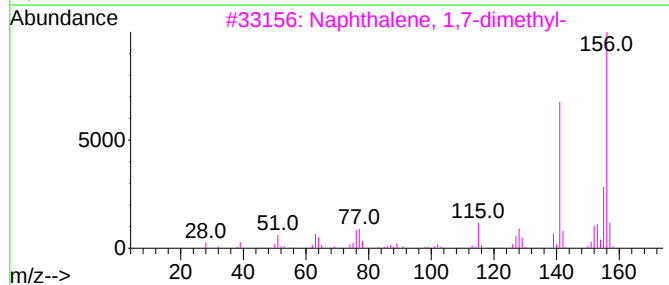
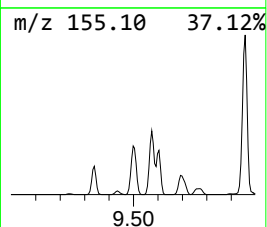
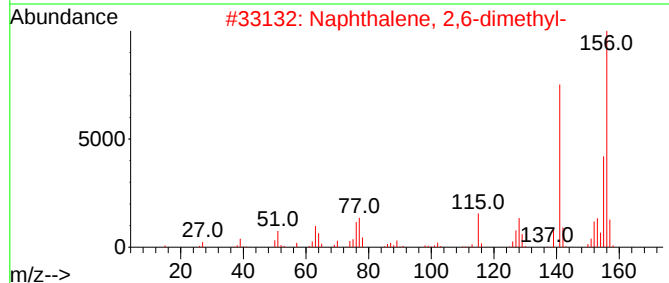
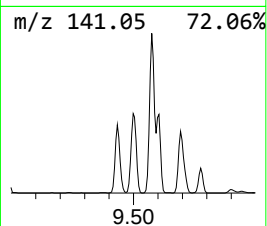
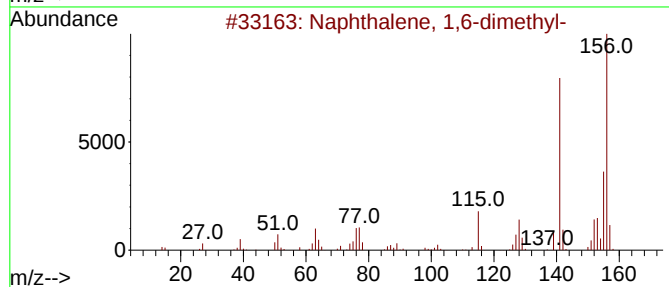
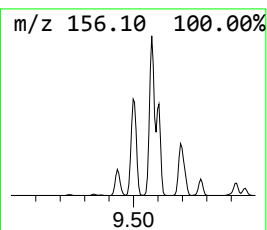
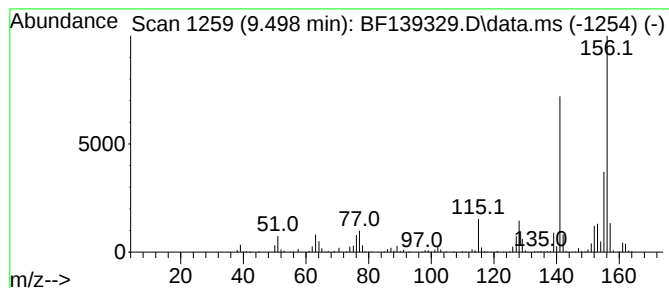
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Naphthalene, 1,6-dimethyl- Concentration Rank 3

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.498 | 7.89 ng | 464943 | Acenaphthene-d10 | 9.922 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 2    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 3    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 4    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |
| 5    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
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Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

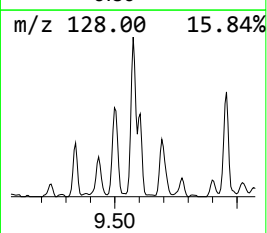
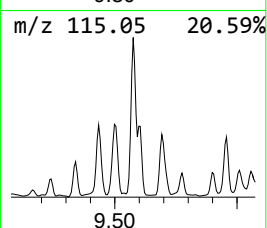
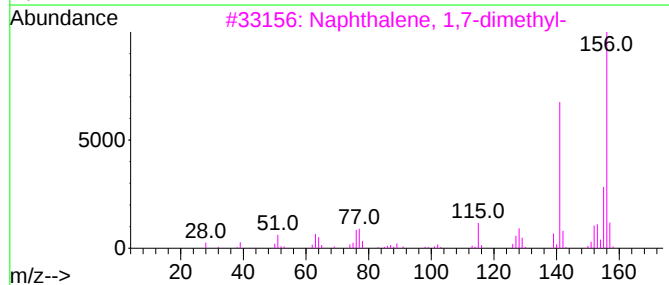
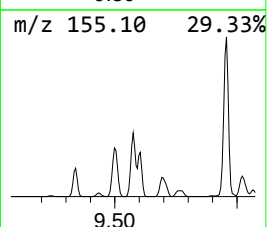
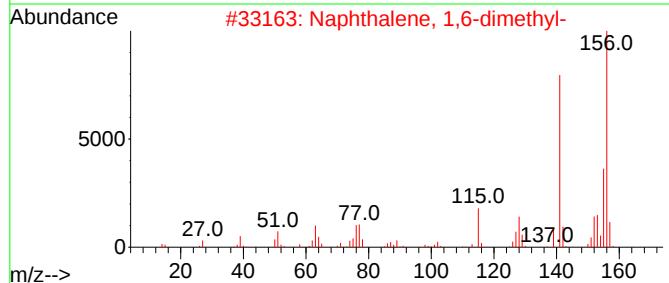
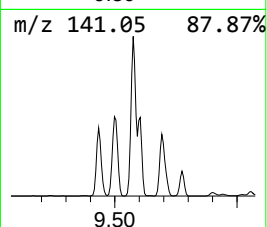
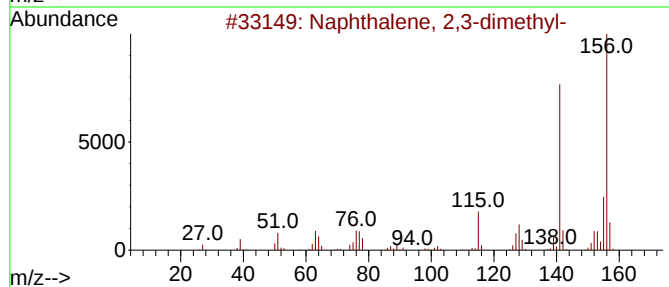
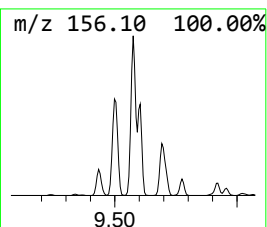
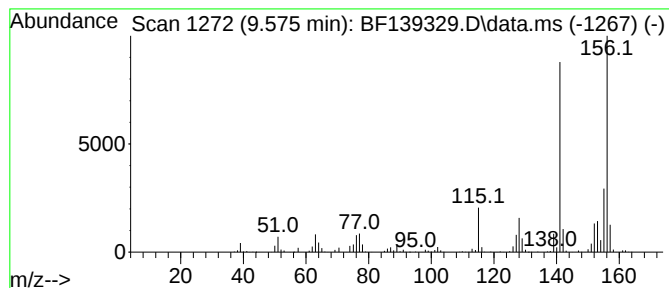
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BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.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 8 Naphthalene, 2,3-dimethyl- Concentration Rank 2

| R.T.      | EstConc                    | Area   | Relative to ISTD | R.T.        |      |
|-----------|----------------------------|--------|------------------|-------------|------|
| 9.575     | 12.08 ng                   | 711783 | Acenaphthene-d10 | 9.922       |      |
| Hit# of 5 | Tentative ID               | MW     | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2,3-dimethyl- | 156    | C12H12           | 000581-40-8 | 97   |
| 2         | Naphthalene, 1,6-dimethyl- | 156    | C12H12           | 000575-43-9 | 97   |
| 3         | Naphthalene, 1,7-dimethyl- | 156    | C12H12           | 000575-37-1 | 97   |
| 4         | Naphthalene, 1,3-dimethyl- | 156    | C12H12           | 000575-41-7 | 97   |
| 5         | Naphthalene, 1,4-dimethyl- | 156    | C12H12           | 000571-58-4 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

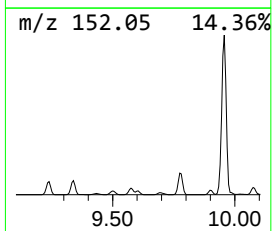
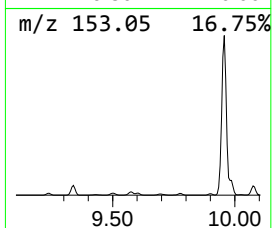
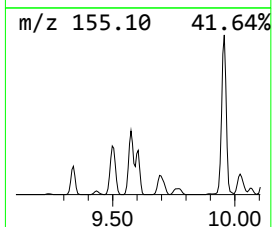
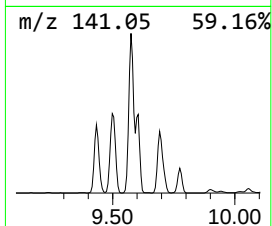
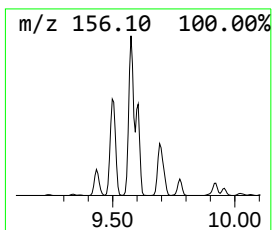
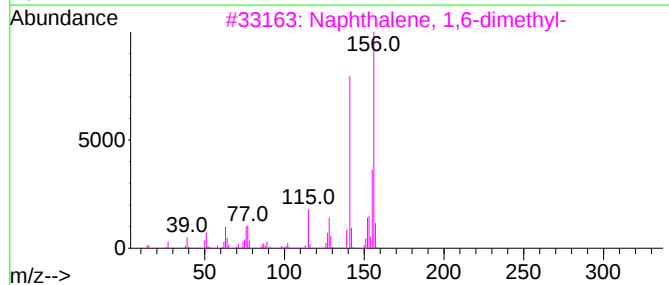
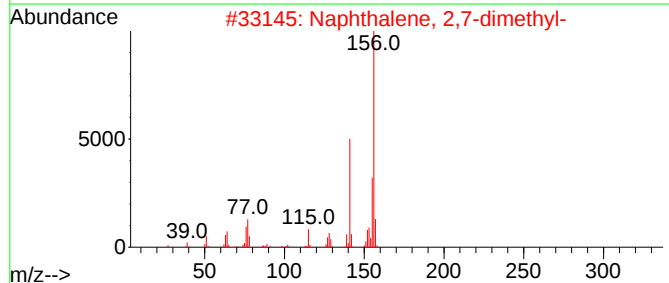
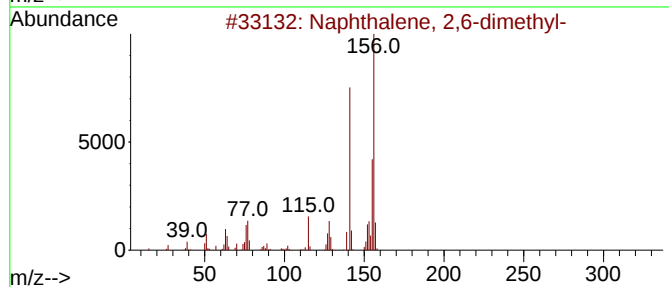
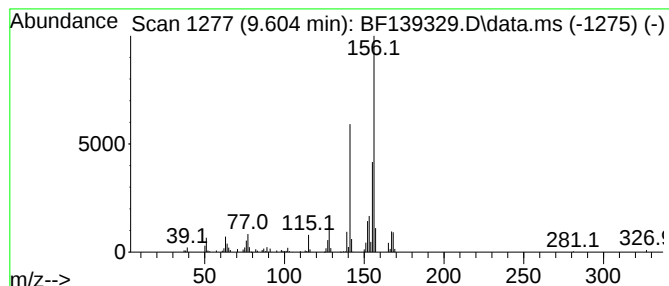
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Naphthalene, 2,6-dimethyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.604 | 4.75 ng | 279662 | Acenaphthene-d10 | 9.922 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 91   |
| 2    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 91   |
| 3    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 90   |
| 4    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 76   |
| 5    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

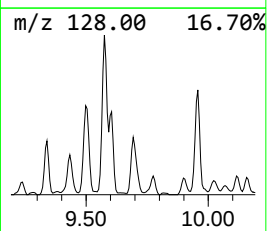
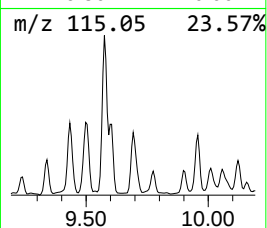
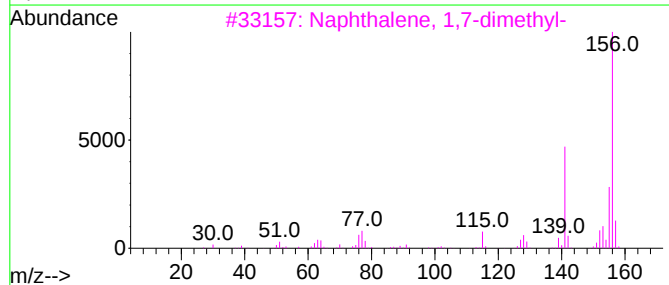
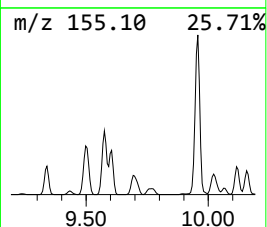
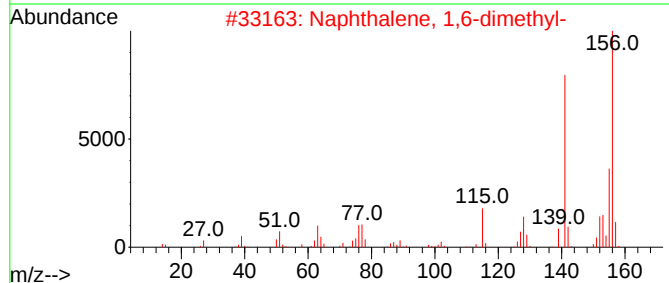
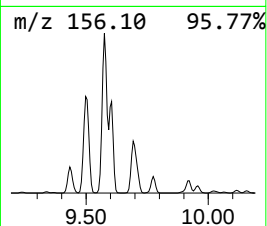
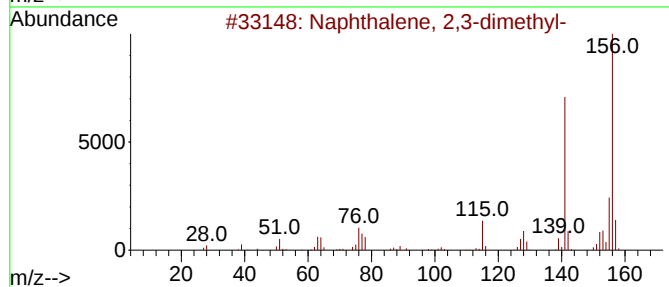
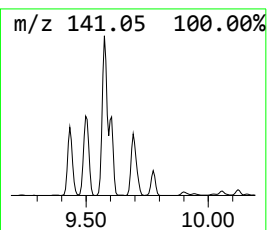
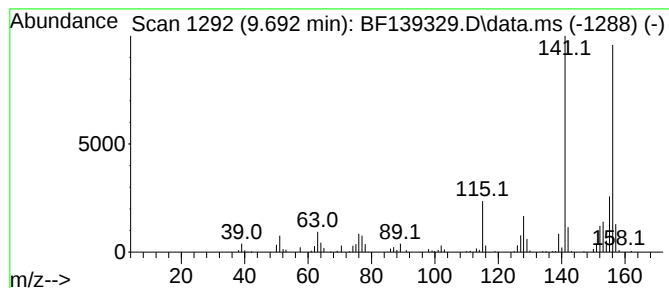
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Naphthalene, 1,7-dimethyl- Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.692 | 4.67 ng | 275285 | Acenaphthene-d10 | 9.922 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 2    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 3    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |
| 4    |    |   | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 96   |
| 5    |    |   | Naphthalene, 1,8-dimethyl- | 156 | C12H12  | 000569-41-5 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

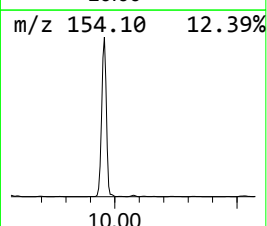
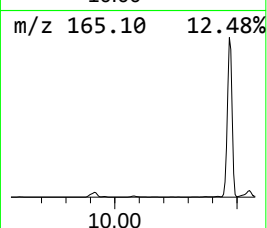
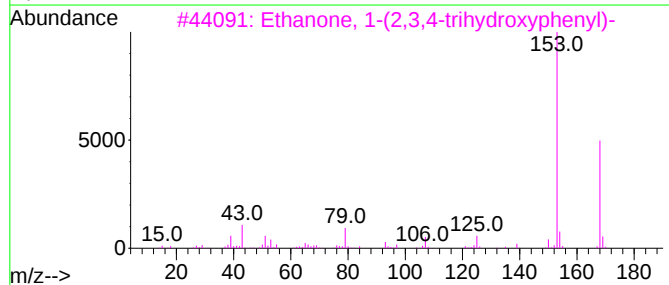
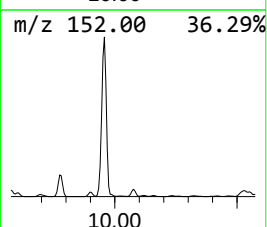
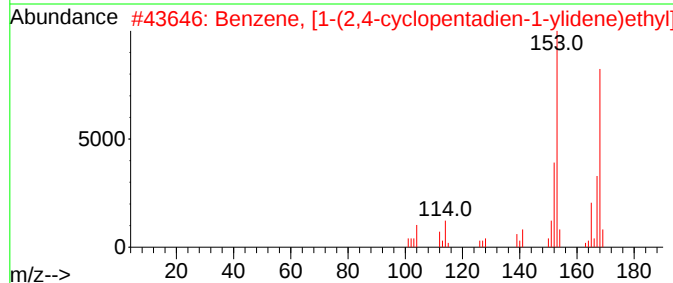
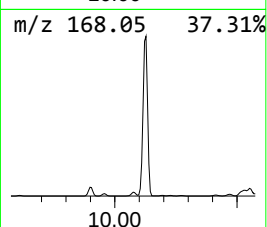
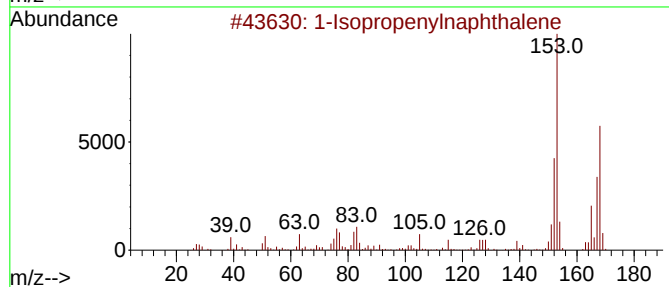
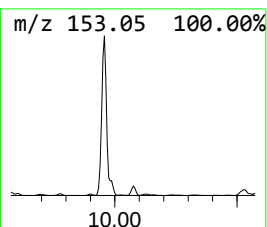
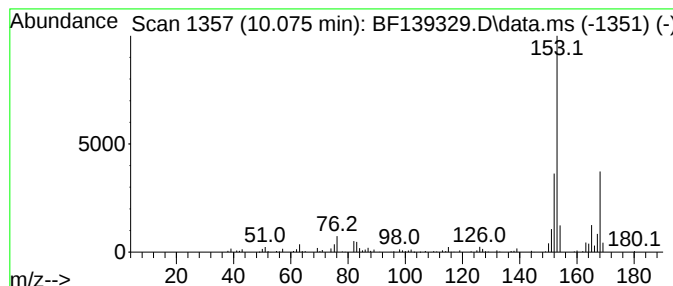
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 1-Isopropenylnaphthalene Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 10.075    | 3.98 ng                             | 234687 | Acenaphthene-d10 | 9.922       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 1-Isopropenylnaphthalene            | 168    | C13H12           | 001855-47-6 | 83   |
| 2         | Benzene, [1-(2,4-cyclopentadien-... | 168    | C13H12           | 002320-32-3 | 72   |
| 3         | Ethanone, 1-(2,3,4-trihydroxyphe... | 168    | C8H8O4           | 000528-21-2 | 58   |
| 4         | Thieno[2,3-b]thiophene, 2-ethyl-    | 168    | C8H8S2           | 005912-01-6 | 53   |
| 5         | Ethanone, 1-(2,4,6-trihydroxyphe... | 168    | C8H8O4           | 000480-66-0 | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

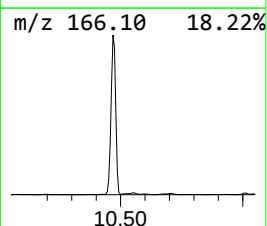
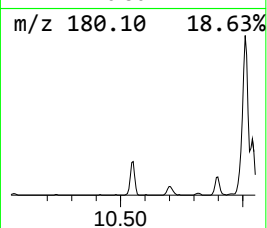
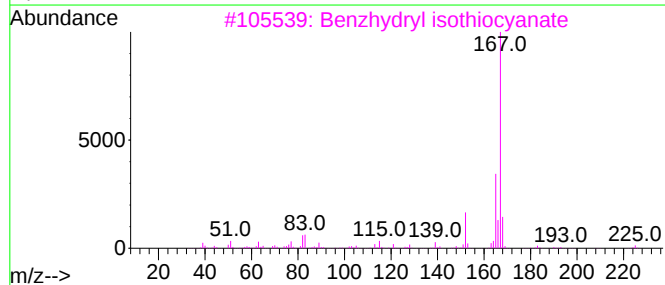
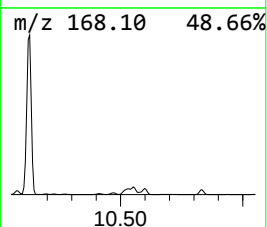
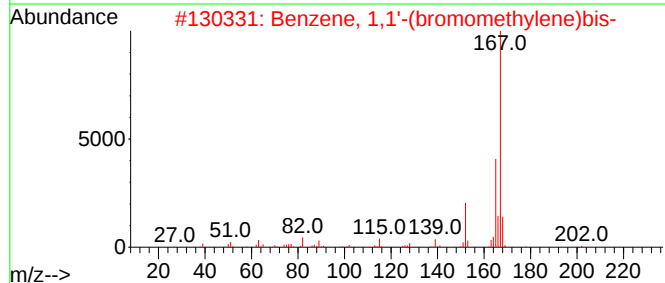
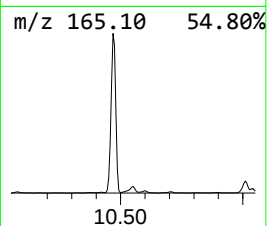
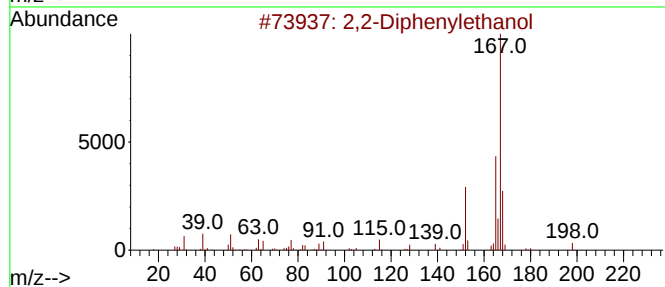
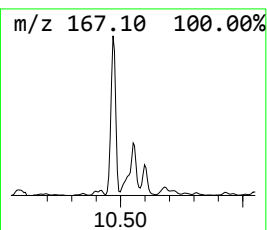
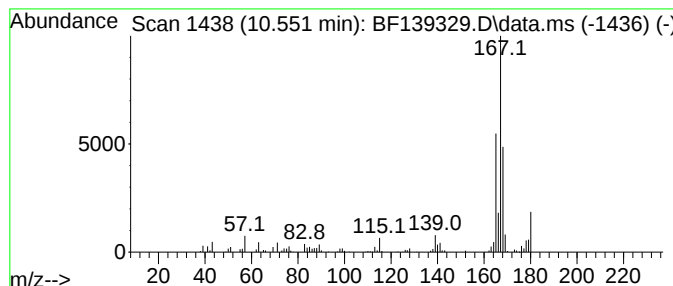
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 2,2-Diphenylethanol Concentration Rank 15

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.  |
|--------|---------|--------|------------------|-------|
| 10.551 | 2.63 ng | 155082 | Acenaphthene-d10 | 9.922 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2,2-Diphenylethanol                 | 198 | C14H14O  | 001883-32-5 | 59   |
| 2    |    |   | Benzene, 1,1'-(bromomethylene)bis-  | 246 | C13H11Br | 000776-74-9 | 53   |
| 3    |    |   | Benzhydryl isothiocyanate           | 225 | C14H11NS | 003550-21-8 | 47   |
| 4    |    |   | 2-Propanone, 1,1-diphenyl-          | 210 | C15H14O  | 000781-35-1 | 47   |
| 5    |    |   | Benzeneacetaldehyde, .alpha.-phe... | 196 | C14H12O  | 000947-91-1 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

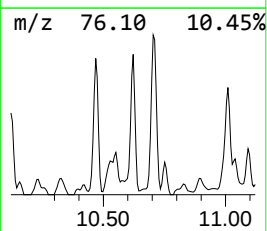
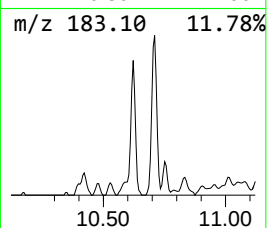
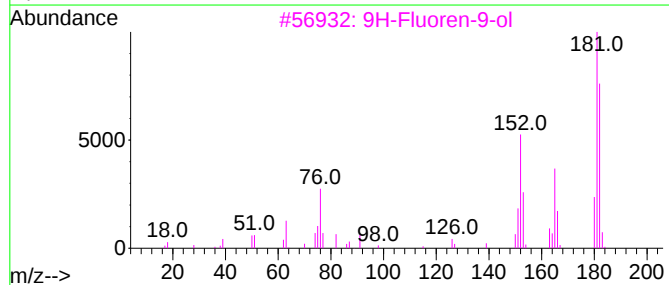
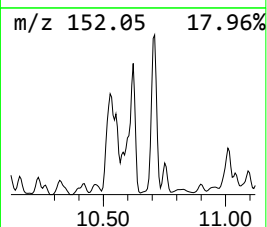
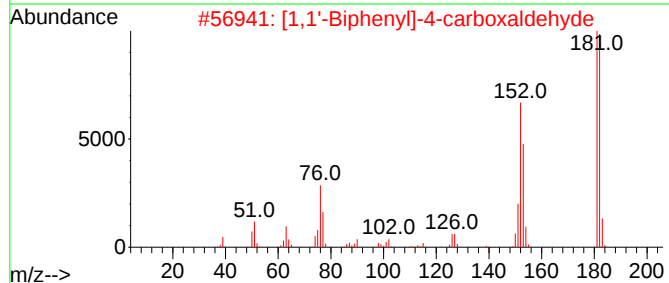
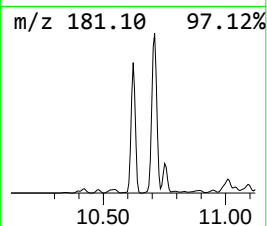
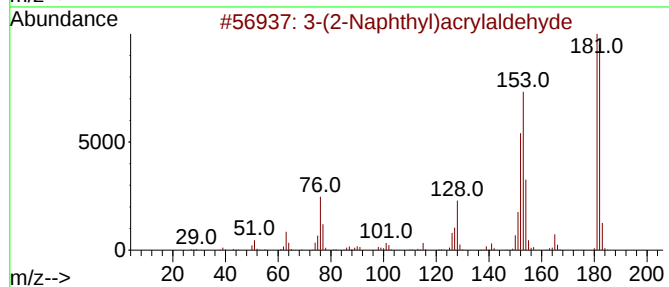
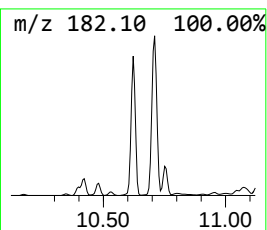
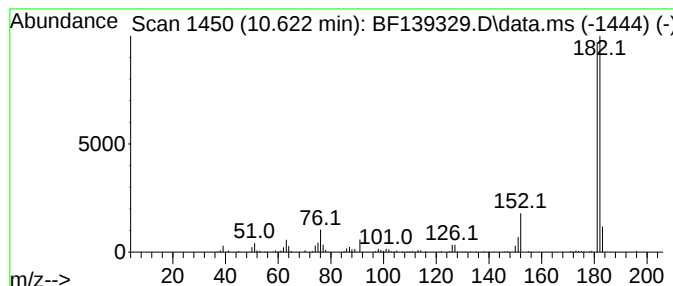
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 3-(2-Naphthyl)acrylaldehyde Concentration Rank 4

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.  |
|--------|---------|--------|------------------|-------|
| 10.622 | 7.47 ng | 440013 | Acenaphthene-d10 | 9.922 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 3-(2-Naphthyl)acrylaldehyde         | 182 | C13H10O | 020884-05-3 | 86   |
| 2    |      | [1,1'-Biphenyl]-4-carboxaldehyde    | 182 | C13H10O | 003218-36-8 | 83   |
| 3    |      | 9H-Fluoren-9-ol                     | 182 | C13H10O | 001689-64-1 | 78   |
| 4    |      | 9H-Xanthene                         | 182 | C13H10O | 000092-83-1 | 72   |
| 5    |      | 2,4,6-Cycloheptatrien-1-one, 2-p... | 182 | C13H10O | 014562-09-5 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
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Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

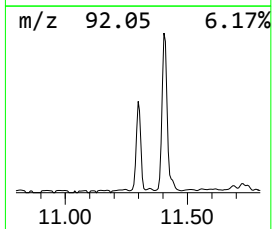
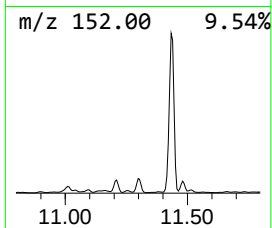
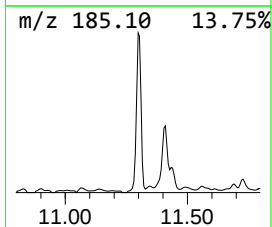
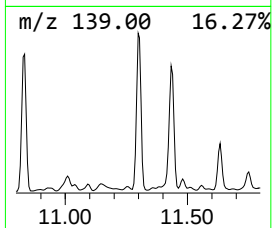
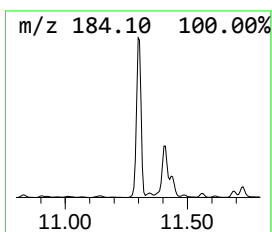
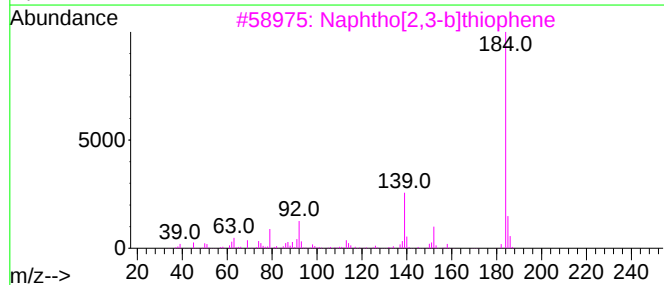
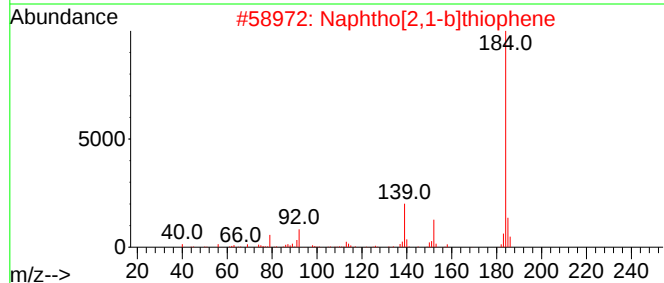
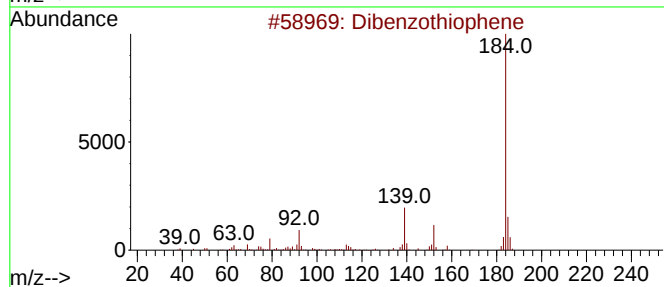
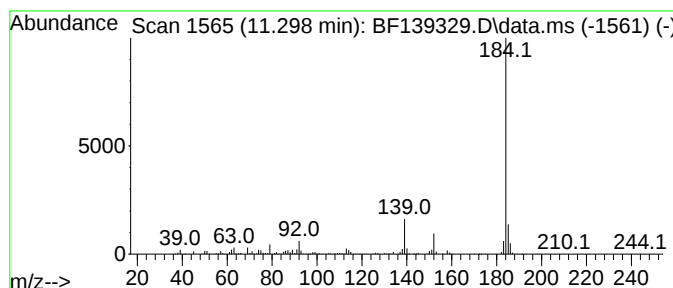
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ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.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 Dibenzothiophene Concentration Rank 18

| R.T.      | EstConc                   | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|---------------------------|--------|------------------|---------|-------------|------|
| 11.298    | 2.01 ng                   | 431805 | Phenanthrene-d10 |         | 11.410      |      |
| Hit# of 5 | Tentative ID              |        | MW               | MolForm | CAS#        | Qual |
| 1         | Dibenzothiophene          |        | 184              | C12H8S  | 000132-65-0 | 96   |
| 2         | Naphtho[2,1-b]thiophene   |        | 184              | C12H8S  | 000233-02-3 | 95   |
| 3         | Naphtho[2,3-b]thiophene   |        | 184              | C12H8S  | 000268-77-9 | 93   |
| 4         | Azuleno(2,1-b)thiophene   |        | 184              | C12H8S  | 000248-13-5 | 91   |
| 5         | 1,1'-Biphenyl, 2-methoxy- |        | 184              | C13H12O | 000086-26-0 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

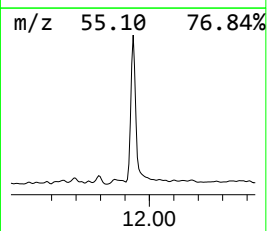
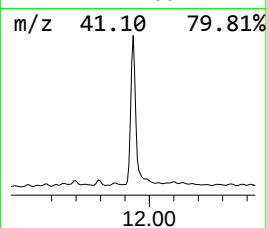
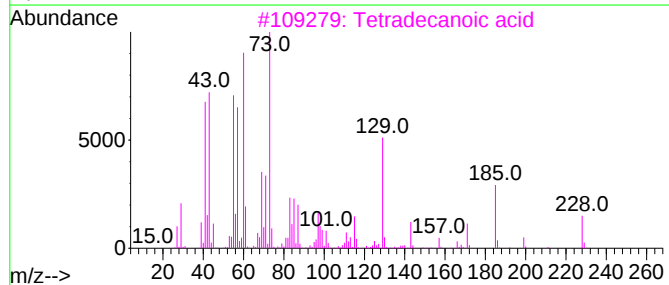
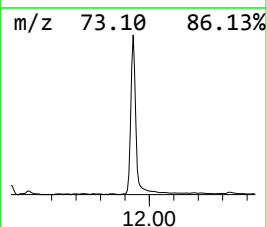
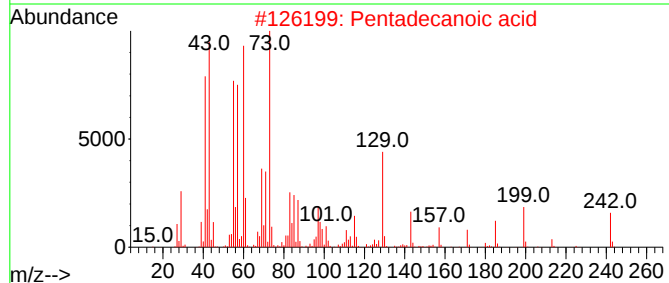
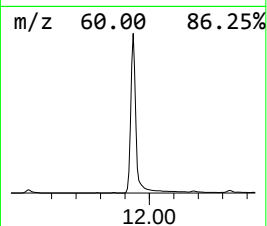
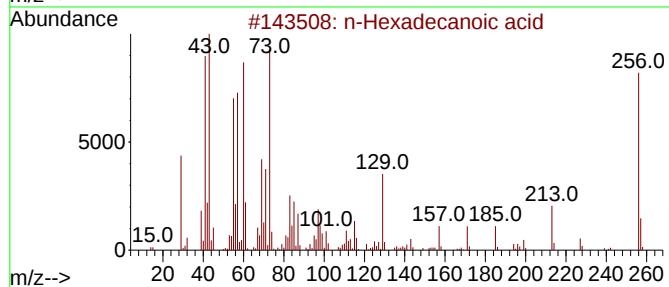
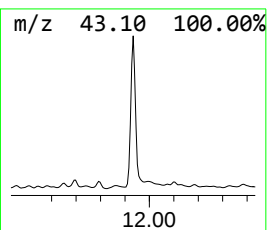
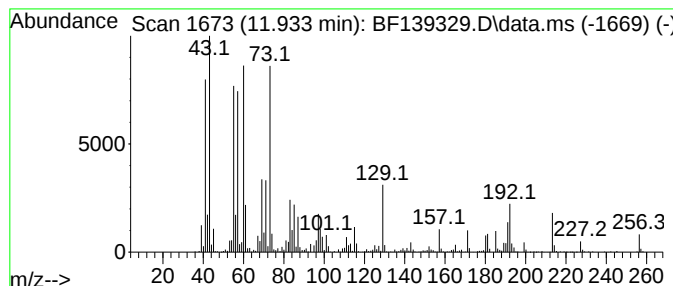
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 n-Hexadecanoic acid Concentration Rank 6

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 11.933 | 5.36 ng | 1151370 | Phenanthrene-d10 | 11.410 |

| Hit# | of 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|------|---------------------|-----|----------|-------------|------|
| 1    |      | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 2    |      | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 76   |
| 3    |      | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 74   |
| 4    |      | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 70   |
| 5    |      | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 60   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

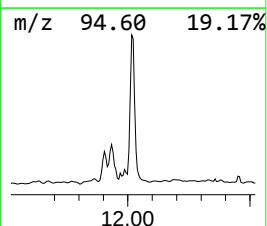
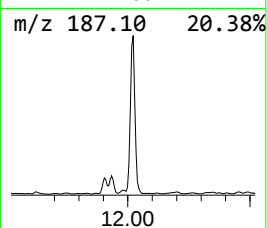
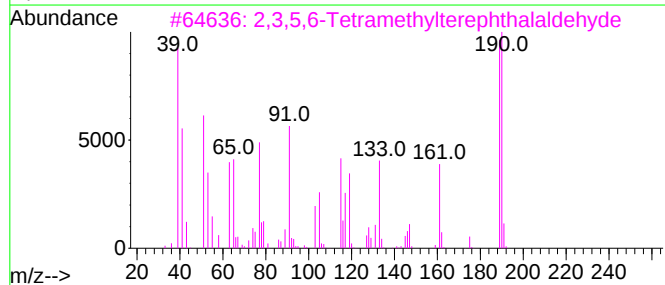
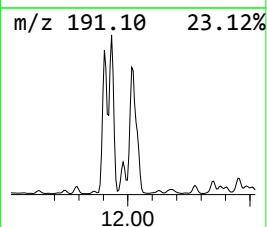
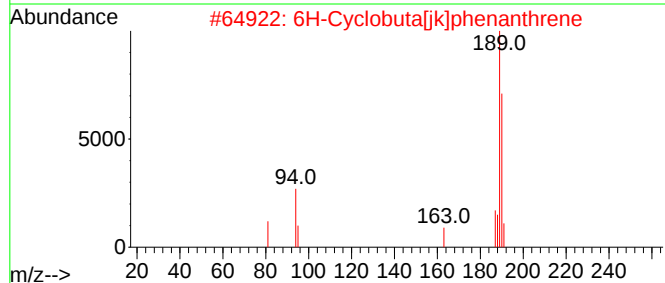
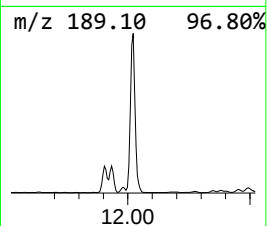
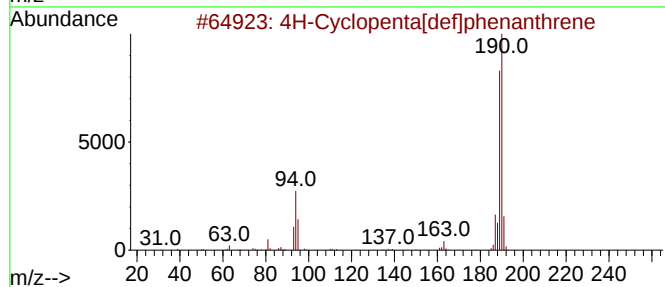
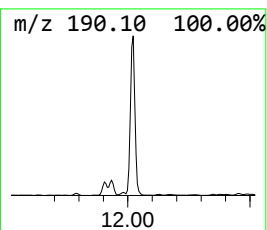
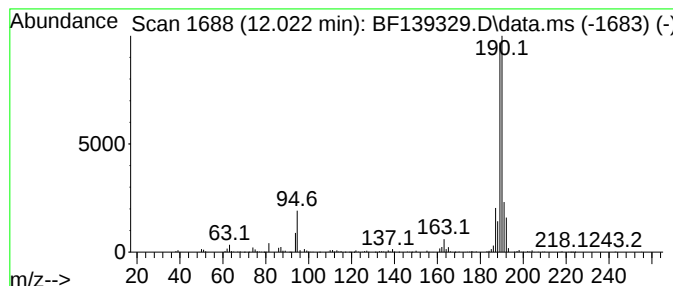
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 4H-Cyclopenta[def]phenanthrene Concentration Rank 13

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 12.022 | 3.29 ng | 707077 | Phenanthrene-d10 | 11.410 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10     | 000203-64-5 | 97   |
| 2    |    |   | 6H-Cyclobuta[jk]phenanthrene        | 190 | C15H10     | 083469-43-6 | 72   |
| 3    |    |   | 2,3,5,6-Tetramethylterephthalald... | 190 | C12H14O2   | 007072-01-7 | 50   |
| 4    |    |   | 2,2'-Bis(4,5-dimethylimidazole)     | 190 | C10H14N4   | 069286-06-2 | 45   |
| 5    |    |   | Thiophene-3-carboxaldehyde, 5-ch... | 190 | C5H4BClO3S | 036155-87-0 | 40   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

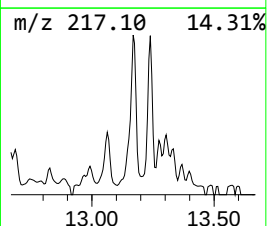
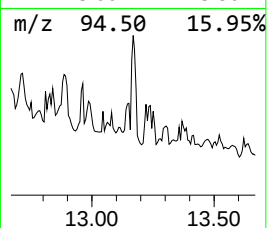
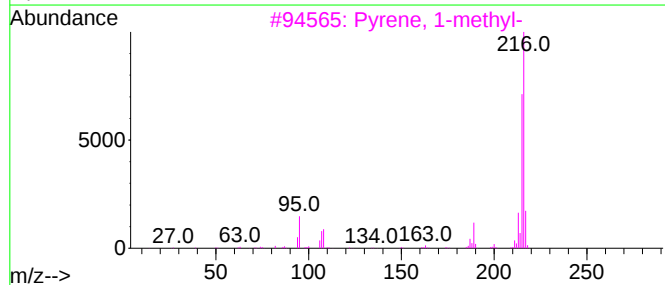
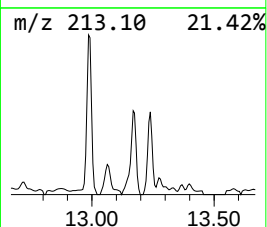
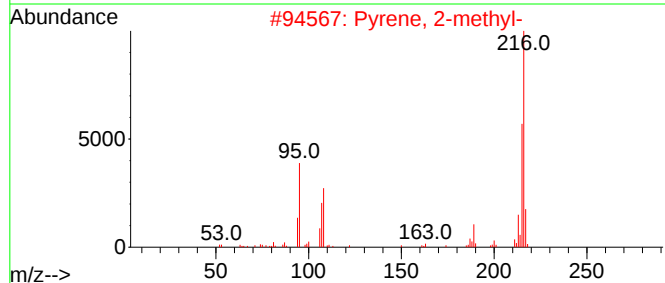
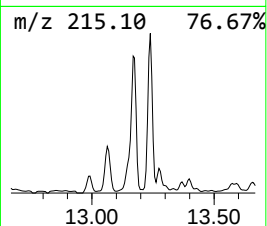
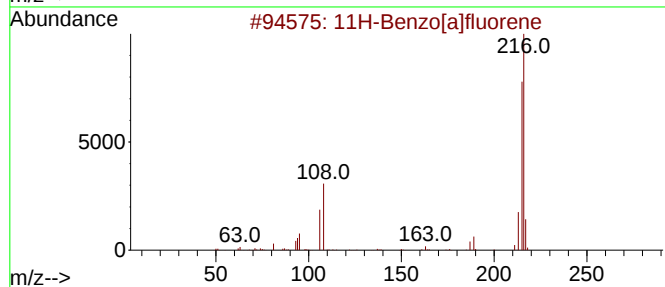
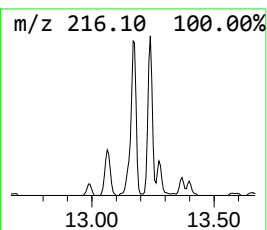
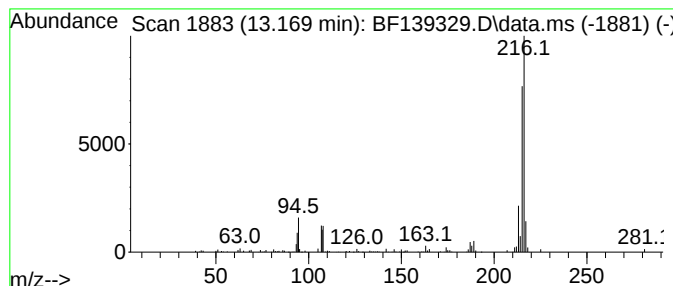
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 17 11H-Benzo[a]fluorene Concentration Rank 16

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 13.169 | 2.37 ng | 77622 | Chrysene-d12     | 14.039 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 89   |
| 2    |    |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 87   |
| 3    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 81   |
| 4    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 74   |
| 5    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

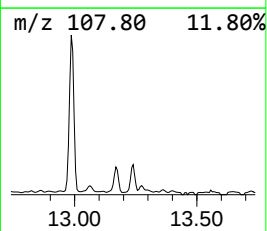
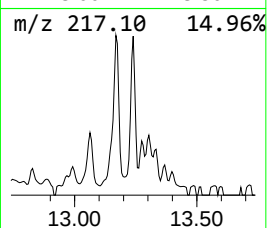
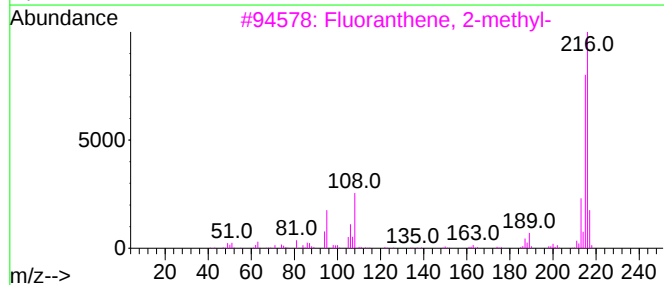
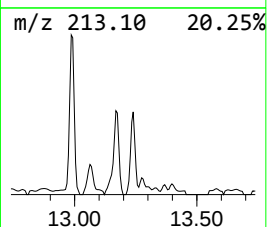
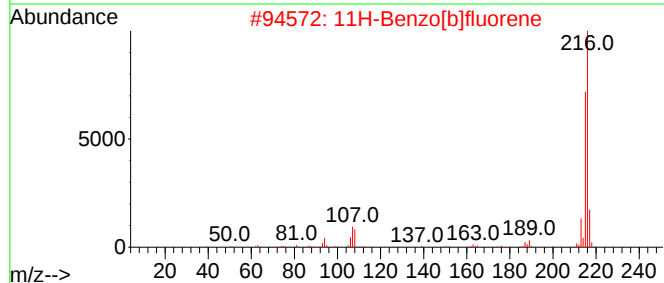
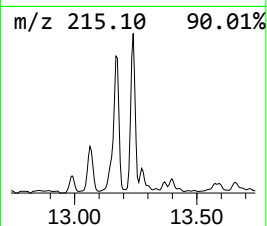
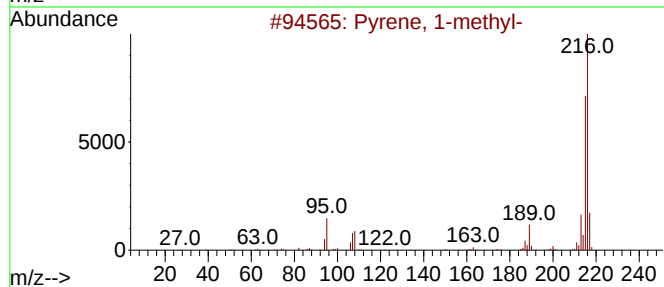
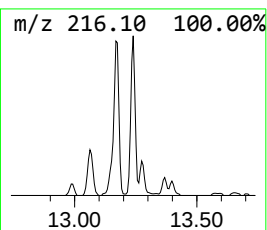
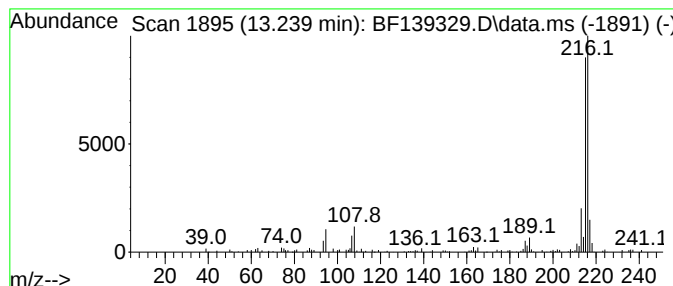
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 18 Pyrene, 1-methyl- Concentration Rank 17

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 13.239 | 2.21 ng | 72447 | Chrysene-d12     | 14.039 |

| 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 | 93   |
| 3    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 91   |
| 4    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 87   |
| 5    |    |   | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

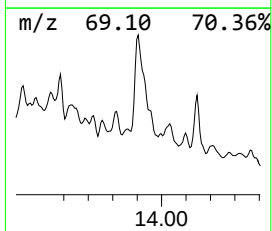
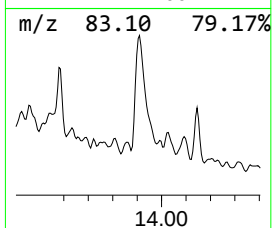
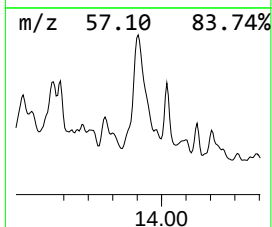
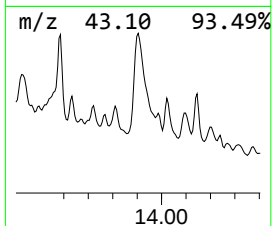
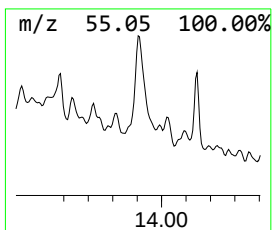
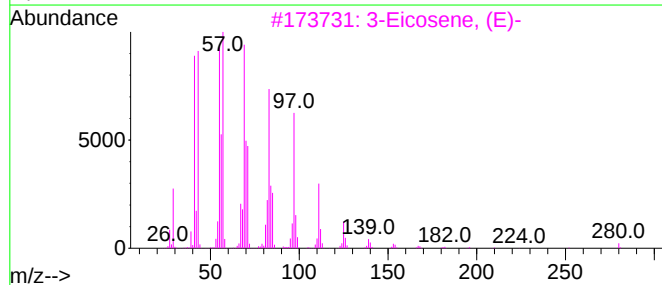
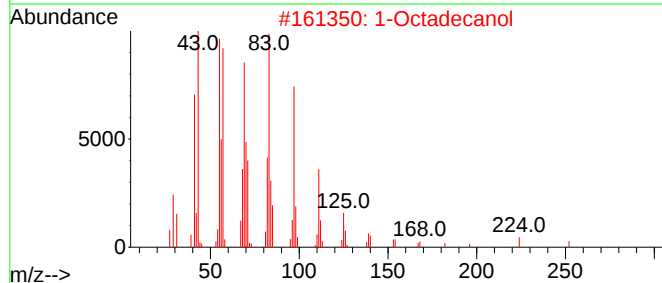
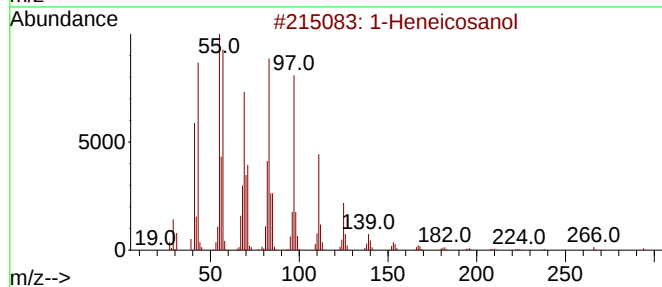
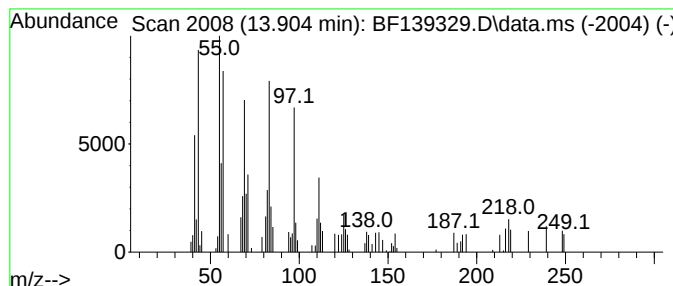
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 1-Heneicosanol Concentration Rank 19

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 13.904 | 2.01 ng | 65778 | Chrysene-d12     | 14.039 |

| Hit# | of 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------|-----|---------|-------------|------|
| 1    |      | 1-Heneicosanol      | 312 | C21H44O | 015594-90-8 | 90   |
| 2    |      | 1-Octadecanol       | 270 | C18H38O | 000112-92-5 | 74   |
| 3    |      | 3-Eicosene, (E)-    | 280 | C20H40  | 074685-33-9 | 72   |
| 4    |      | 17-Pentatriacontene | 491 | C35H70  | 006971-40-0 | 72   |
| 5    |      | 1-Hexacosanol       | 382 | C26H54O | 000506-52-5 | 62   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091124\  
Data File : BF139329.D  
Acq On : 11 Sep 2024 16:50  
Operator : RC/JU  
Sample : P3912-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
282-COMPOSITE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF090424.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 |
| Butane, 2-metho... | 2.181  | 132.2   | ng    | 4231510  | 1                     | 6.881  | 640020  | 20.0 |
| 1,3-Dioxolane, ... | 2.999  | 7.1     | ng    | 226199   | 1                     | 6.881  | 640020  | 20.0 |
| 3-Penten-2-one,... | 4.475  | 5.0     | ng    | 158543   | 1                     | 6.881  | 640020  | 20.0 |
| 2-Pentanone, 4-... | 5.093  | 2.8     | ng    | 89579    | 1                     | 6.881  | 640020  | 20.0 |
| 1-Phenoxypropan... | 8.475  | 5.0     | ng    | 332513   | 2                     | 8.163  | 1343530 | 20.0 |
| Naphthalene, 1-... | 9.434  | 3.4     | ng    | 202338   | 3                     | 9.922  | 1178060 | 20.0 |
| Naphthalene, 1,... | 9.498  | 7.9     | ng    | 464943   | 3                     | 9.922  | 1178060 | 20.0 |
| Naphthalene, 2,... | 9.575  | 12.1    | ng    | 711783   | 3                     | 9.922  | 1178060 | 20.0 |
| Naphthalene, 2,... | 9.604  | 4.8     | ng    | 279662   | 3                     | 9.922  | 1178060 | 20.0 |
| Naphthalene, 1,... | 9.692  | 4.7     | ng    | 275285   | 3                     | 9.922  | 1178060 | 20.0 |
| 1-Isopropenylna... | 10.075 | 4.0     | ng    | 234687   | 3                     | 9.922  | 1178060 | 20.0 |
| 2,2-Diphenyleth... | 10.551 | 2.6     | ng    | 155082   | 3                     | 9.922  | 1178060 | 20.0 |
| 3-(2-Naphthyl)a... | 10.622 | 7.5     | ng    | 440013   | 3                     | 9.922  | 1178060 | 20.0 |
| Dibenzothiophene   | 11.298 | 2.0     | ng    | 431805   | 4                     | 11.410 | 4299460 | 20.0 |
| n-Hexadecanoic ... | 11.933 | 5.4     | ng    | 1151370  | 4                     | 11.410 | 4299460 | 20.0 |
| 4H-Cyclopenta[d... | 12.022 | 3.3     | ng    | 707077   | 4                     | 11.410 | 4299460 | 20.0 |
| 11H-Benzo[a]flu... | 13.169 | 2.4     | ng    | 77622    | 5                     | 14.039 | 655417  | 20.0 |
| Pyrene, 1-methyl-  | 13.239 | 2.2     | ng    | 72447    | 5                     | 14.039 | 655417  | 20.0 |
| 1-Heneicosanol     | 13.904 | 2.0     | ng    | 65778    | 5                     | 14.039 | 655417  | 20.0 |