

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF090120\  
 Data File : BF121508.D  
 Acq On : 1 Sep 2020 21:13  
 Operator : JU/CG  
 Sample : L3506-11 5X  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OR-02-083120

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 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-BF082720.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| 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         | 5.463       | 570           | 574         | 582          | rVB      | 898545         | 884891        | 5.76%           | 1.033%        |
| 2         | 6.475       | 742           | 746         | 754          | rBV      | 1019712        | 918863        | 5.98%           | 1.073%        |
| 3         | 6.857       | 807           | 811         | 814          | rBV      | 1874662        | 1563710       | 10.18%          | 1.826%        |
| 4         | 7.410       | 902           | 905         | 911          | rVB      | 624462         | 580647        | 3.78%           | 0.678%        |
| 5         | 8.139       | 1024          | 1029        | 1031         | rBV      | 2435858        | 2069398       | 13.47%          | 2.416%        |
| 6         | 8.163       | 1031          | 1033        | 1036         | rVB      | 548140         | 417220        | 2.72%           | 0.487%        |
| 7         | 8.851       | 1148          | 1150        | 1154         | rVB      | 298062         | 259679        | 1.69%           | 0.303%        |
| 8         | 8.951       | 1163          | 1167        | 1172         | rBV2     | 263663         | 310309        | 2.02%           | 0.362%        |
| 9         | 9.216       | 1208          | 1212        | 1215         | rBV      | 1660747        | 1239687       | 8.07%           | 1.448%        |
| 10        | 9.298       | 1223          | 1226        | 1228         | rBV      | 214021         | 205077        | 1.33%           | 0.239%        |
| 11        | 9.480       | 1253          | 1257        | 1261         | rVB2     | 213851         | 271392        | 1.77%           | 0.317%        |
| 12        | 9.557       | 1261          | 1270        | 1272         | rBV      | 305392         | 365603        | 2.38%           | 0.427%        |
| 13        | 9.580       | 1272          | 1274        | 1276         | rVV      | 216961         | 179239        | 1.17%           | 0.209%        |
| 14        | 9.610       | 1276          | 1279        | 1283         | rVV2     | 268933         | 352583        | 2.30%           | 0.412%        |
| 15        | 9.675       | 1287          | 1290        | 1294         | rVB2     | 171227         | 189150        | 1.23%           | 0.221%        |
| 16        | 9.757       | 1301          | 1304        | 1309         | rVB      | 423295         | 350457        | 2.28%           | 0.409%        |
| 17        | 9.828       | 1314          | 1316        | 1321         | rVB      | 307503         | 258442        | 1.68%           | 0.302%        |
| 18        | 9.898       | 1321          | 1328        | 1331         | rBV      | 2706814        | 2314595       | 15.07%          | 2.703%        |
| 19        | 9.928       | 1331          | 1333        | 1337         | rVB2     | 238398         | 225868        | 1.47%           | 0.264%        |
| 20        | 10.098      | 1359          | 1362        | 1367         | rVB      | 365699         | 381546        | 2.48%           | 0.446%        |
| 21        | 10.139      | 1367          | 1369        | 1375         | rBV3     | 114574         | 158022        | 1.03%           | 0.185%        |
| 22        | 10.327      | 1399          | 1401        | 1404         | rVB      | 403144         | 323414        | 2.11%           | 0.378%        |
| 23        | 10.445      | 1417          | 1421        | 1427         | rBV      | 600884         | 540599        | 3.52%           | 0.631%        |
| 24        | 10.551      | 1434          | 1439        | 1445         | rBV      | 192459         | 272791        | 1.78%           | 0.319%        |
| 25        | 10.680      | 1457          | 1461        | 1466         | rBV      | 755347         | 777095        | 5.06%           | 0.907%        |
| 26        | 10.804      | 1479          | 1482        | 1490         | rBV      | 571945         | 657548        | 4.28%           | 0.768%        |
| 27        | 10.910      | 1497          | 1500        | 1503         | rVB      | 281228         | 211732        | 1.38%           | 0.247%        |
| 28        | 10.992      | 1510          | 1514        | 1517         | rBV2     | 169781         | 209770        | 1.37%           | 0.245%        |
| 29        | 11.186      | 1544          | 1547        | 1552         | rVB      | 245720         | 228267        | 1.49%           | 0.267%        |
| 30        | 11.251      | 1556          | 1558        | 1561         | rVV      | 486702         | 436202        | 2.84%           | 0.509%        |
| 31        | 11.280      | 1561          | 1563        | 1569         | rVB2     | 419496         | 432306        | 2.81%           | 0.505%        |
| 32        | 11.386      | 1577          | 1581        | 1583         | rBV      | 2344600        | 2034750       | 13.24%          | 2.376%        |
| 33        | 11.410      | 1583          | 1585        | 1589         | rVV      | 3324903        | 2600354       | 16.93%          | 3.036%        |
| 34        | 11.463      | 1589          | 1594        | 1597         | rVB      | 623967         | 581496        | 3.79%           | 0.679%        |

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 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OR-02-083120

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BF082720.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |          |          |         |         |
|----|--------|------|------|------|------|----------|----------|---------|---------|
| 35 | 11.616 | 1614 | 1620 | 1622 | rBV  | 268382   | 324296   | 2.11%   | 0.379%  |
| 36 | 11.680 | 1628 | 1631 | 1633 | rBV  | 522762   | 421697   | 2.74%   | 0.492%  |
| 37 | 11.704 | 1633 | 1635 | 1644 | rVB2 | 623922   | 615597   | 4.01%   | 0.719%  |
| 38 | 11.786 | 1644 | 1649 | 1652 | rBV  | 8212104  | 7549763  | 49.14%  | 8.816%  |
| 39 | 11.886 | 1663 | 1666 | 1668 | rBV  | 508980   | 397842   | 2.59%   | 0.465%  |
| 40 | 11.916 | 1668 | 1671 | 1674 | rVB  | 933065   | 756952   | 4.93%   | 0.884%  |
| 41 | 11.998 | 1682 | 1685 | 1688 | rBV  | 653540   | 680143   | 4.43%   | 0.794%  |
| 42 | 12.022 | 1688 | 1689 | 1693 | rVB  | 365248   | 226248   | 1.47%   | 0.264%  |
| 43 | 12.092 | 1698 | 1701 | 1709 | rVB2 | 375218   | 515710   | 3.36%   | 0.602%  |
| 44 | 12.186 | 1714 | 1717 | 1719 | rVV  | 432247   | 397311   | 2.59%   | 0.464%  |
| 45 | 12.204 | 1719 | 1720 | 1724 | rVB2 | 298511   | 253367   | 1.65%   | 0.296%  |
| 46 | 12.439 | 1757 | 1760 | 1762 | rBV  | 388563   | 378478   | 2.46%   | 0.442%  |
| 47 | 12.480 | 1762 | 1767 | 1768 | rVV2 | 9323367  | 11979299 | 77.98%  | 13.988% |
| 48 | 12.504 | 1768 | 1771 | 1777 | rVV  | 11748581 | 15362462 | 100.00% | 17.939% |
| 49 | 12.580 | 1780 | 1784 | 1786 | rBV  | 4649815  | 4446759  | 28.95%  | 5.193%  |
| 50 | 12.604 | 1786 | 1788 | 1791 | rVB  | 2872975  | 2179888  | 14.19%  | 2.545%  |
| 51 | 12.698 | 1801 | 1804 | 1806 | rVB  | 175108   | 172702   | 1.12%   | 0.202%  |
| 52 | 12.833 | 1824 | 1827 | 1829 | rVB  | 2091910  | 1712578  | 11.15%  | 2.000%  |
| 53 | 12.974 | 1848 | 1851 | 1854 | rVV  | 1235221  | 992959   | 6.46%   | 1.159%  |
| 54 | 13.057 | 1860 | 1865 | 1867 | rBV2 | 207453   | 343562   | 2.24%   | 0.401%  |
| 55 | 13.139 | 1875 | 1879 | 1880 | rBV3 | 410713   | 407474   | 2.65%   | 0.476%  |
| 56 | 13.157 | 1880 | 1882 | 1886 | rVB  | 467764   | 453802   | 2.95%   | 0.530%  |
| 57 | 13.227 | 1888 | 1894 | 1897 | rBV2 | 715518   | 942708   | 6.14%   | 1.101%  |
| 58 | 13.263 | 1897 | 1900 | 1903 | rVB2 | 251514   | 211467   | 1.38%   | 0.247%  |
| 59 | 13.304 | 1904 | 1907 | 1909 | rBV  | 435635   | 371410   | 2.42%   | 0.434%  |
| 60 | 13.357 | 1914 | 1916 | 1919 | rBV  | 143906   | 193691   | 1.26%   | 0.226%  |
| 61 | 13.551 | 1947 | 1949 | 1952 | rBV3 | 178743   | 157237   | 1.02%   | 0.184%  |
| 62 | 13.663 | 1966 | 1968 | 1973 | rVB  | 230081   | 231588   | 1.51%   | 0.270%  |
| 63 | 13.710 | 1973 | 1976 | 1978 | rBV2 | 261482   | 327063   | 2.13%   | 0.382%  |
| 64 | 13.774 | 1984 | 1987 | 1990 | rBV2 | 203900   | 263847   | 1.72%   | 0.308%  |
| 65 | 13.880 | 2001 | 2005 | 2010 | rVB2 | 420889   | 516187   | 3.36%   | 0.603%  |
| 66 | 13.974 | 2018 | 2021 | 2023 | rVB  | 265589   | 238638   | 1.55%   | 0.279%  |
| 67 | 14.027 | 2025 | 2030 | 2033 | rVV2 | 2401558  | 3139478  | 20.44%  | 3.666%  |
| 68 | 14.057 | 2033 | 2035 | 2042 | rVB  | 1095211  | 935522   | 6.09%   | 1.092%  |
| 69 | 14.415 | 2094 | 2096 | 2099 | rVB  | 245559   | 199351   | 1.30%   | 0.233%  |
| 70 | 15.074 | 2204 | 2208 | 2211 | rBV2 | 829762   | 1248452  | 8.13%   | 1.458%  |
| 71 | 15.380 | 2256 | 2260 | 2266 | rVB  | 483254   | 674694   | 4.39%   | 0.788%  |

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Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

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-BF082720.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |     |         |         |        |        |
|----|--------|------|------|------|-----|---------|---------|--------|--------|
| 72 | 15.439 | 2267 | 2270 | 2277 | rVB | 666210  | 760467  | 4.95%  | 0.888% |
| 73 | 15.504 | 2277 | 2281 | 2285 | rBV | 1391260 | 1853565 | 12.07% | 2.164% |

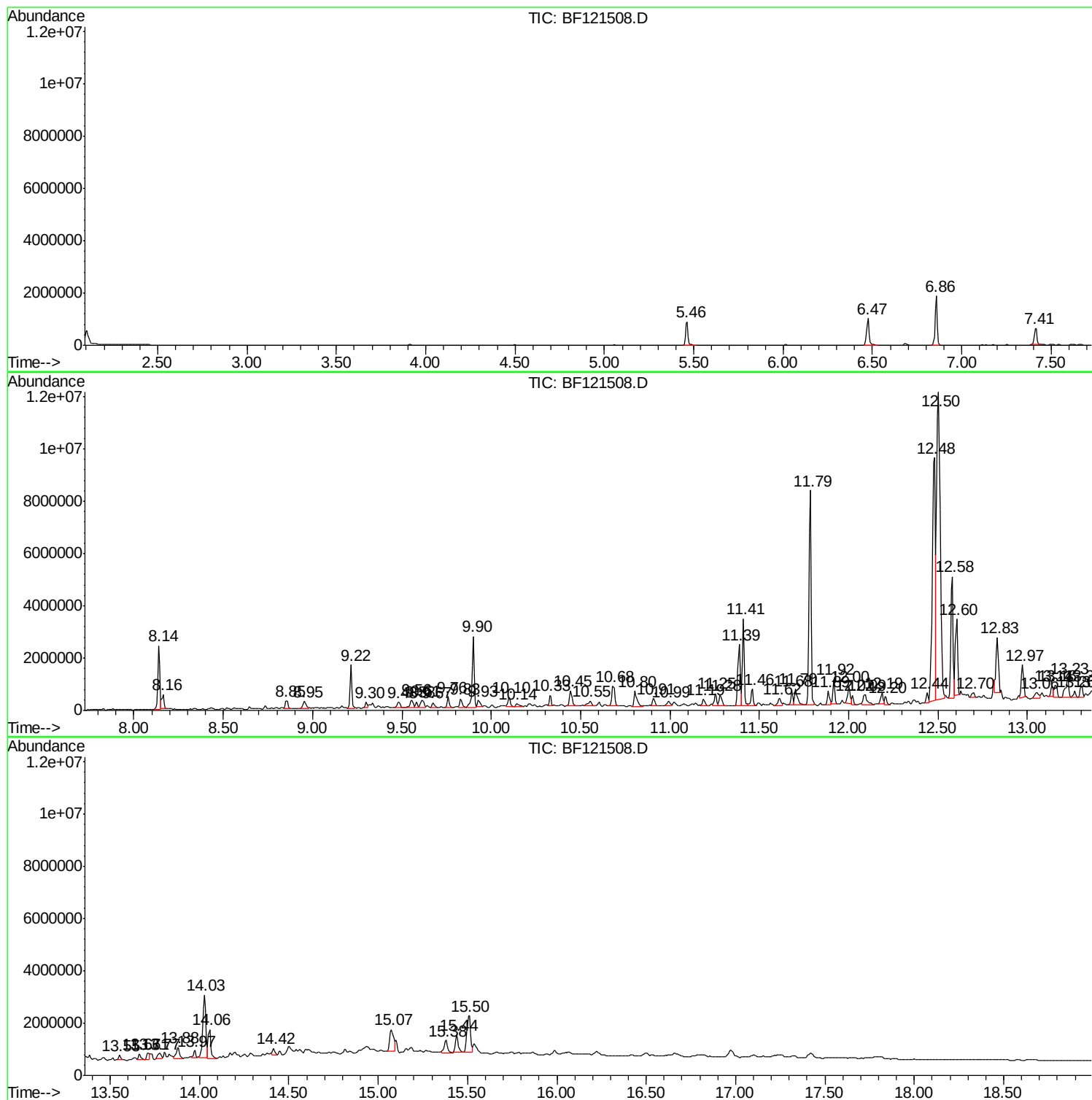
Sum of corrected areas: 85636956

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
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Instrument :  
BNA\_F  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P



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Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

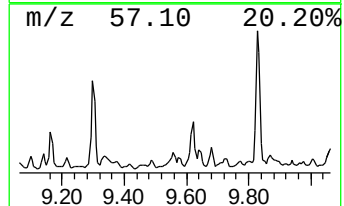
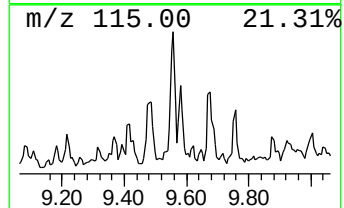
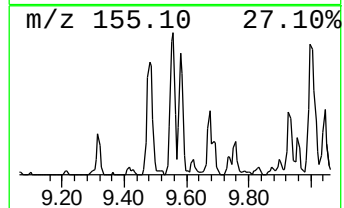
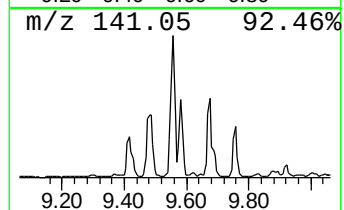
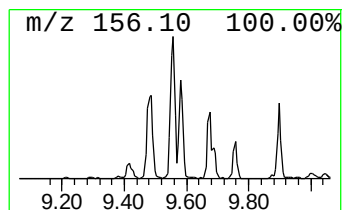
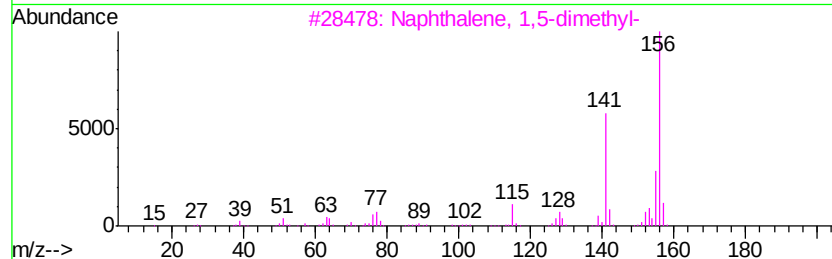
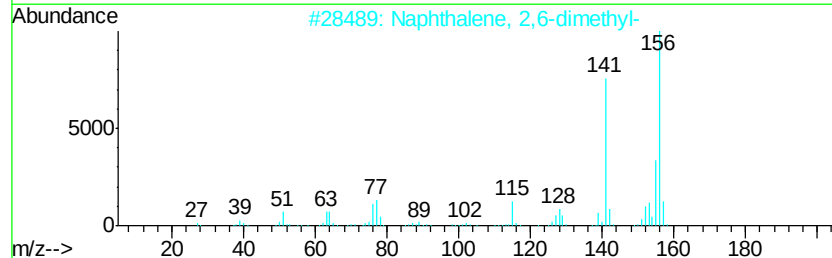
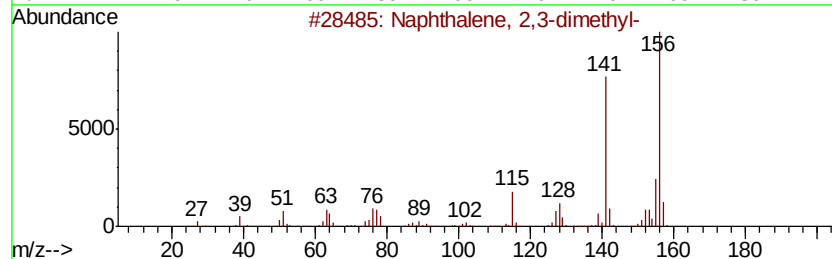
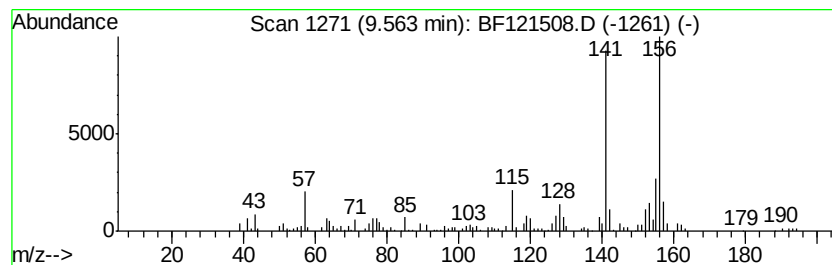
Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Naphthalene, 2,3-dimethyl- Concentration Rank 18

| R.T.      | EstConc                    | Area   | Relative to ISTD |             | R.T. |
|-----------|----------------------------|--------|------------------|-------------|------|
| 9.56      | 3.16 ng                    | 365603 | Acenaphthene-d10 |             | 9.90 |
| Hit# of 5 | Tentative ID               | MW     | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2,3-dimethyl- | 156    | C12H12           | 000581-40-8 | 97   |
| 2         | Naphthalene, 2,6-dimethyl- | 156    | C12H12           | 000581-42-0 | 97   |
| 3         | Naphthalene, 1,5-dimethyl- | 156    | C12H12           | 000571-61-9 | 97   |
| 4         | Naphthalene, 1,6-dimethyl- | 156    | C12H12           | 000575-43-9 | 97   |
| 5         | Naphthalene, 1,3-dimethyl- | 156    | C12H12           | 000575-41-7 | 96   |



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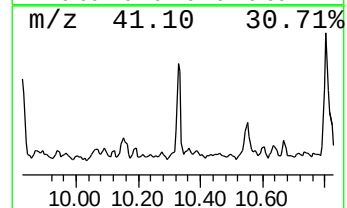
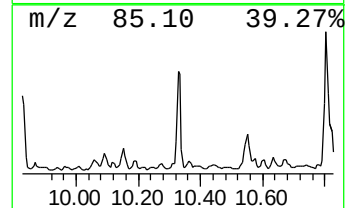
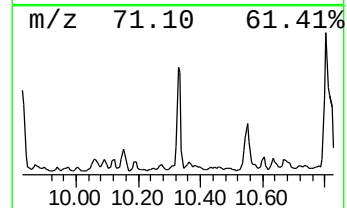
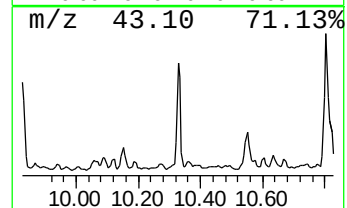
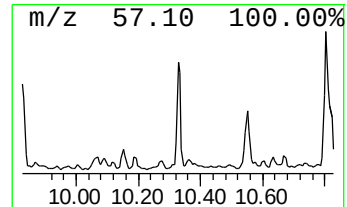
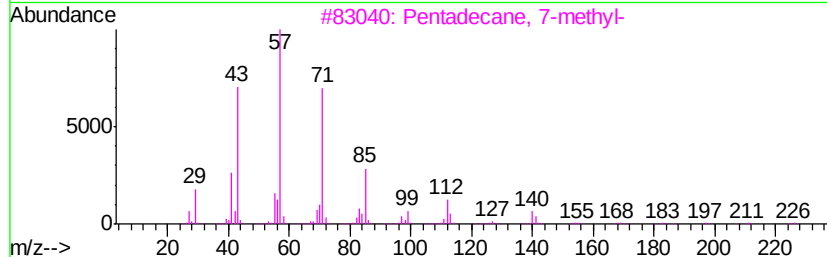
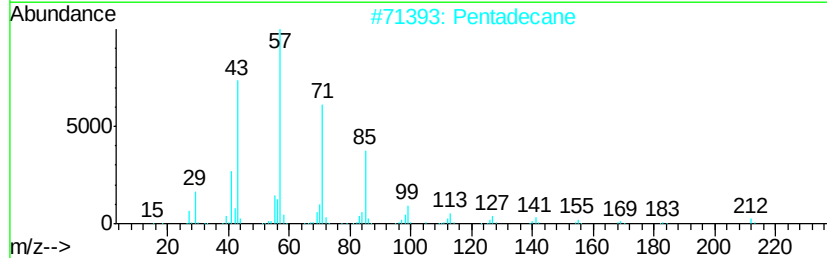
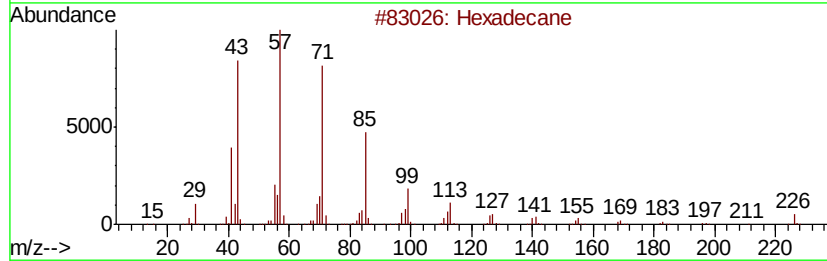
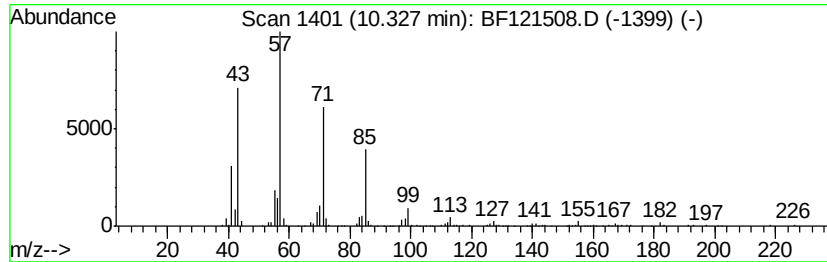
Instrument :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Hexadecane Concentration Rank 20

| R.T.    | EstConc                | Area         | Relative to ISTD | R.T.      |
|---------|------------------------|--------------|------------------|-----------|
| 10.33   | 2.79 ng                | 323414       | Acenaphthene-d10 | 9.90      |
| Hit# of | 5                      | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Hexadecane             | 226 C16H34   | 000544-76-3      | 94        |
| 2       | Pentadecane            | 212 C15H32   | 000629-62-9      | 90        |
| 3       | Pentadecane, 7-methyl- | 226 C16H34   | 006165-40-8      | 87        |
| 4       | Heneicosane            | 296 C21H44   | 000629-94-7      | 83        |
| 5       | Heptadecane            | 240 C17H36   | 000629-78-7      | 83        |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

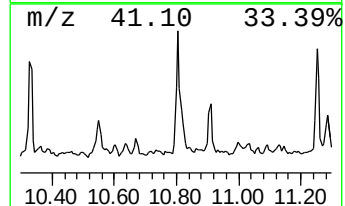
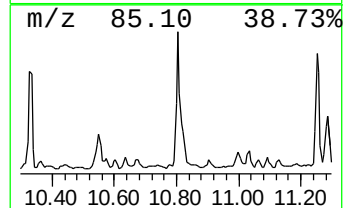
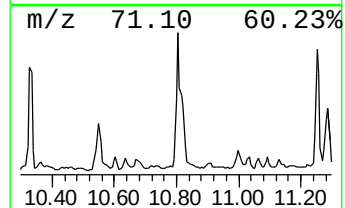
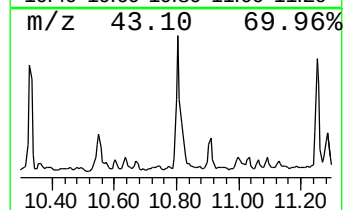
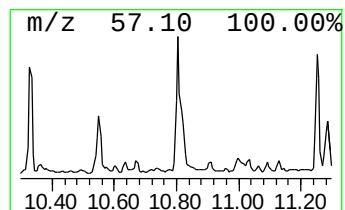
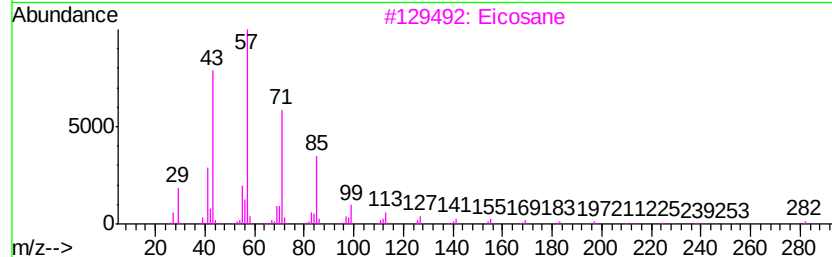
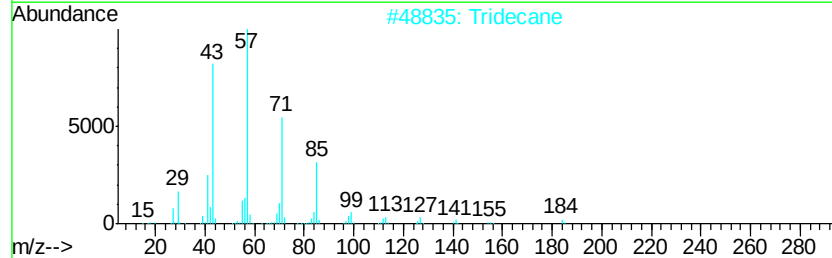
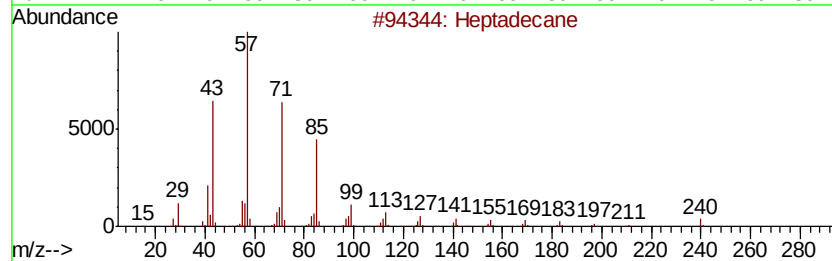
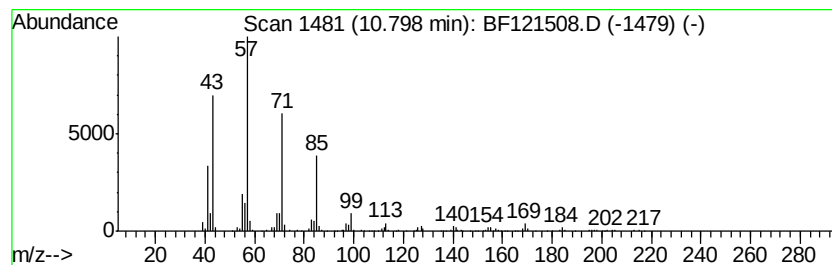
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Heptadecane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.80 | 6.46 ng | 657548 | Phenanthrene-d10 | 11.39 |

| Hit# of | 5                              | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|--------------------------------|--------------|-----|---------|-------------|------|
| 1       | Heptadecane                    |              | 240 | C17H36  | 000629-78-7 | 96   |
| 2       | Tridecane                      |              | 184 | C13H28  | 000629-50-5 | 91   |
| 3       | Eicosane                       |              | 282 | C20H42  | 000112-95-8 | 83   |
| 4       | Tetradecane, 2,6,10-trimethyl- |              | 240 | C17H36  | 014905-56-7 | 80   |
| 5       | Heneicosane                    |              | 296 | C21H44  | 000629-94-7 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

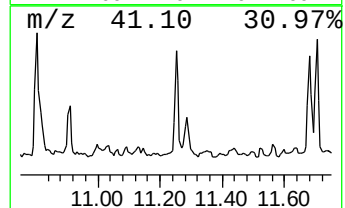
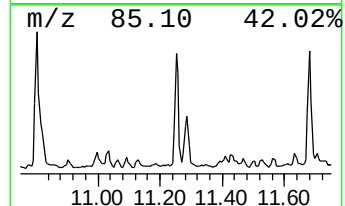
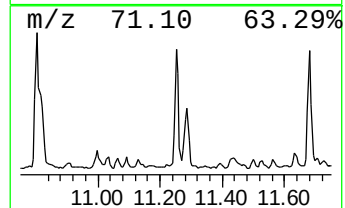
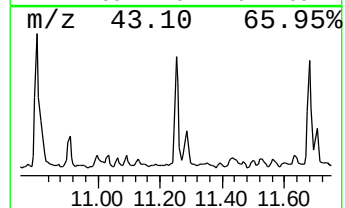
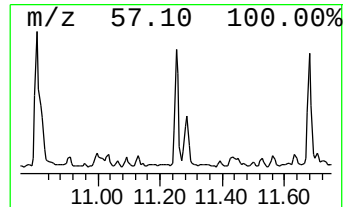
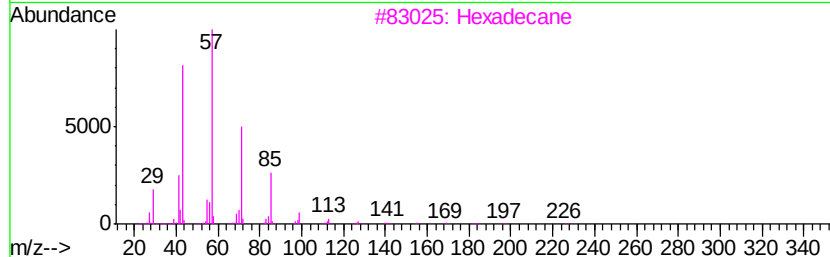
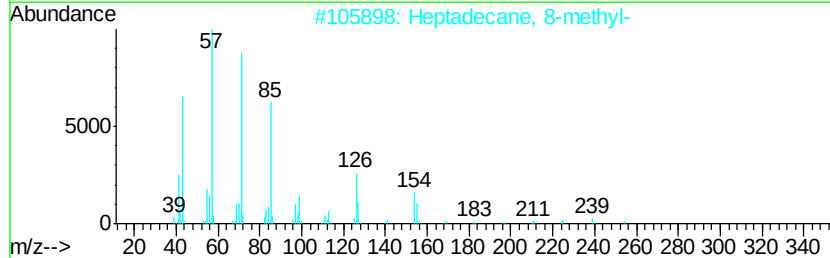
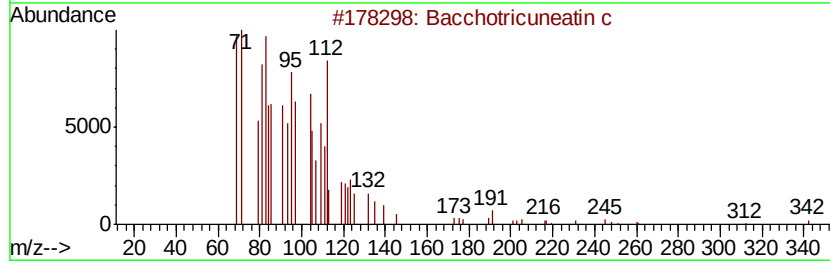
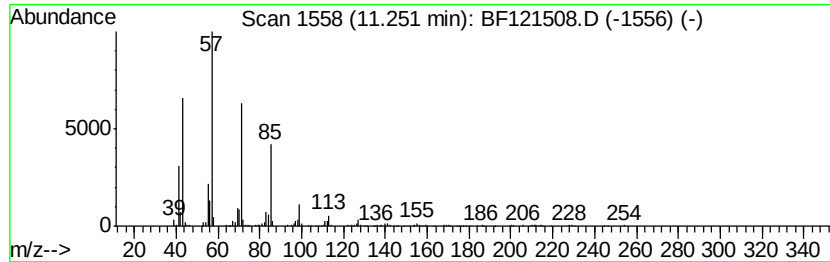
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Bacchotricuneatin c Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.25 | 4.29 ng | 436202 | Phenanthrene-d10 | 11.39 |

| Hit# | of | Tentative ID                 | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------------|-----|----------|-------------|------|
| 1    | 5  | Bacchotricuneatin c          | 342 | C20H22O5 | 066563-30-2 | 97   |
| 2    |    | Heptadecane, 8-methyl-       | 254 | C18H38   | 013287-23-5 | 90   |
| 3    |    | Hexadecane                   | 226 | C16H34   | 000544-76-3 | 90   |
| 4    |    | Tetradecane, 1-iodo-         | 324 | C14H29I  | 019218-94-1 | 90   |
| 5    |    | Dodecane, 2-methyl-6-propyl- | 226 | C16H34   | 055045-08-4 | 90   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
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Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

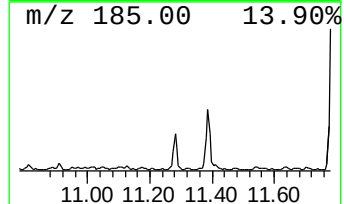
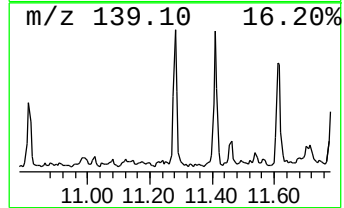
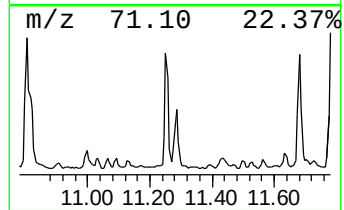
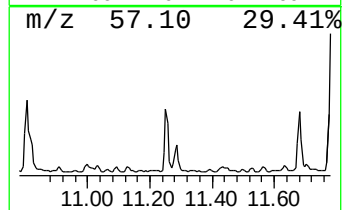
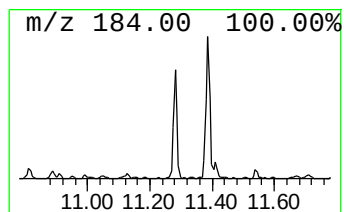
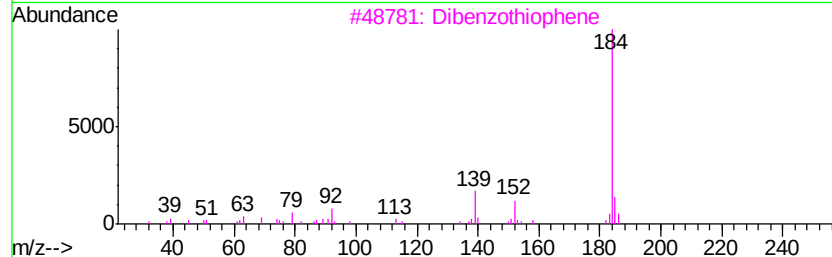
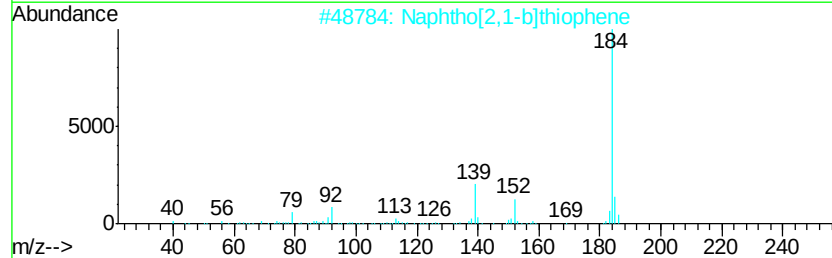
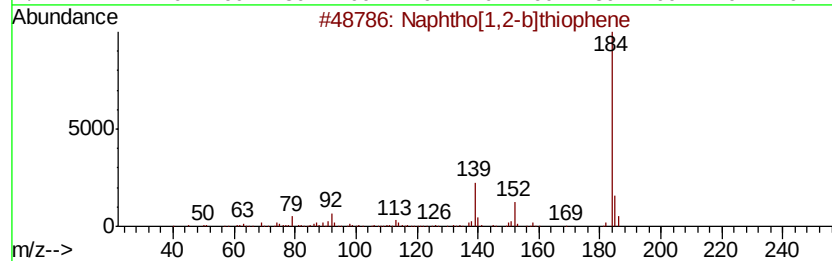
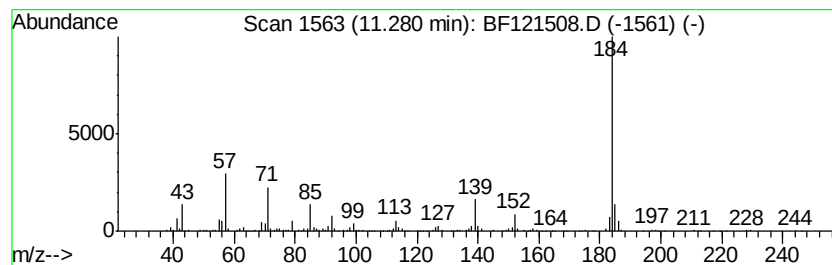
Instrument :  
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ClientSampleId :  
OR-02-083120

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Naphtho[1,2-b]thiophene Concentration Rank 12

| R.T.    | EstConc | Area                    | Relative to ISTD |          | R.T.        |      |
|---------|---------|-------------------------|------------------|----------|-------------|------|
| 11.28   | 4.25 ng | 432306                  | Phenanthrene-d10 |          | 11.39       |      |
| Hit# of | 5       | Tentative ID            | MW               | MolForm  | CAS#        | Qual |
| 1       |         | Naphtho[1,2-b]thiophene | 184              | C12H8S   | 000234-41-3 | 93   |
| 2       |         | Naphtho[2,1-b]thiophene | 184              | C12H8S   | 000233-02-3 | 89   |
| 3       |         | Dibenzothiophene        | 184              | C12H8S   | 000132-65-0 | 89   |
| 4       |         | Azuleno(2,1-b)thiophene | 184              | C12H8S   | 000248-13-5 | 76   |
| 5       |         | Thianthrene, 5-oxide    | 232              | C12H8OS2 | 002362-50-7 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

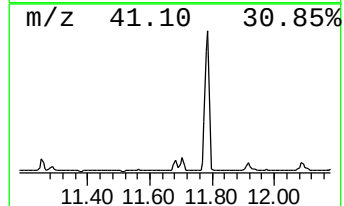
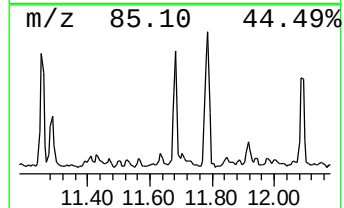
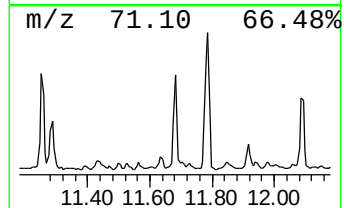
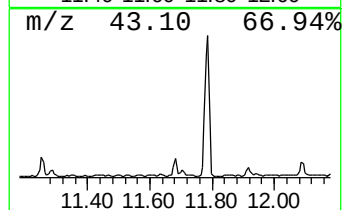
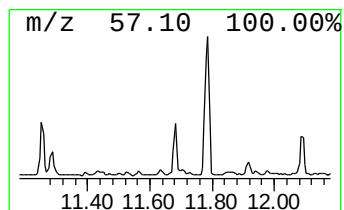
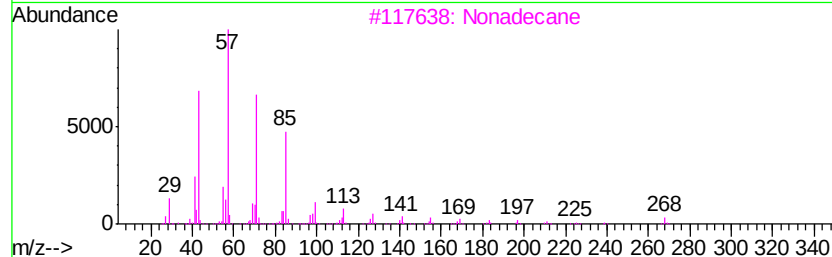
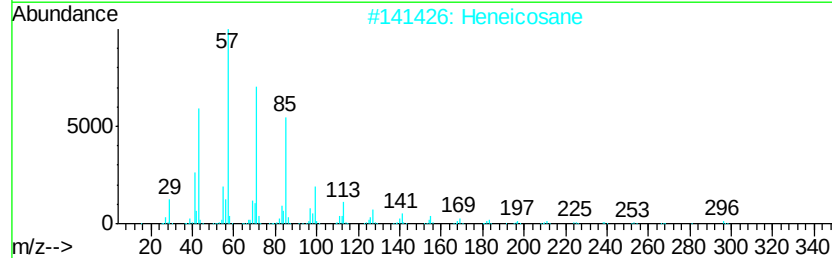
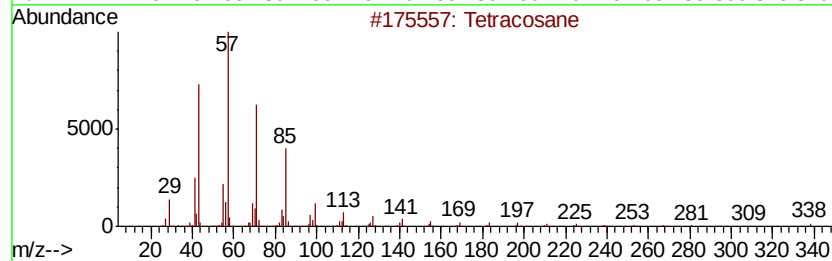
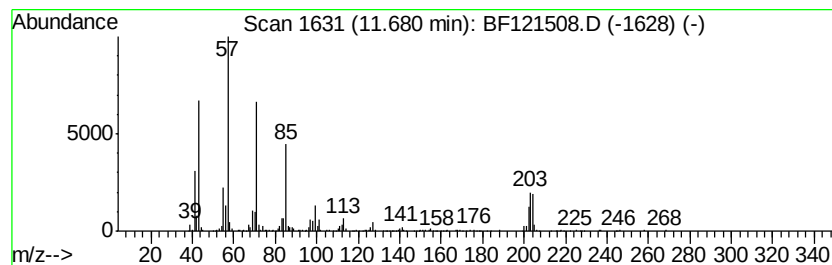
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Tetracosane Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.68 | 4.14 ng | 421697 | Phenanthrene-d10 | 11.39 |

| Hit# of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|---|--------------|-----|---------|-------------|------|
| 1       |   | Tetracosane  | 338 | C24H50  | 000646-31-1 | 62   |
| 2       |   | Heneicosane  | 296 | C21H44  | 000629-94-7 | 58   |
| 3       |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 58   |
| 4       |   | Heptacosane  | 380 | C27H56  | 000593-49-7 | 58   |
| 5       |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
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Sample : L3506-11 5X  
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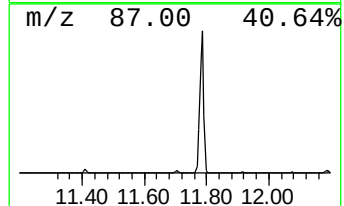
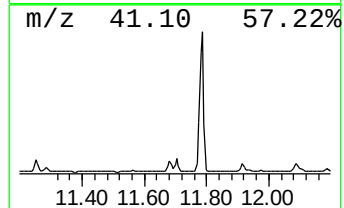
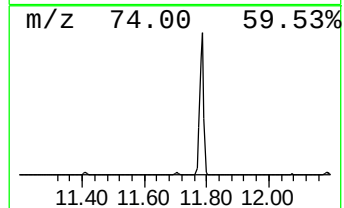
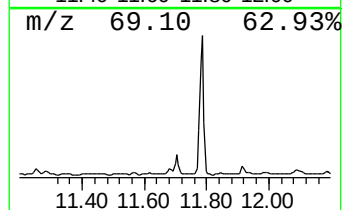
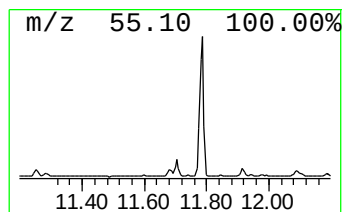
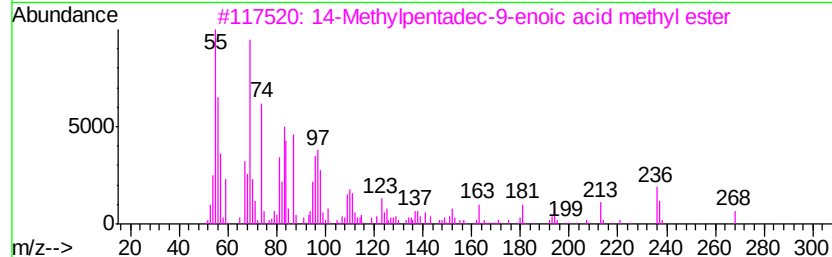
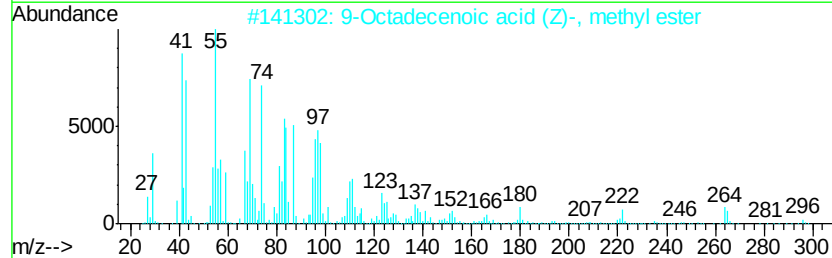
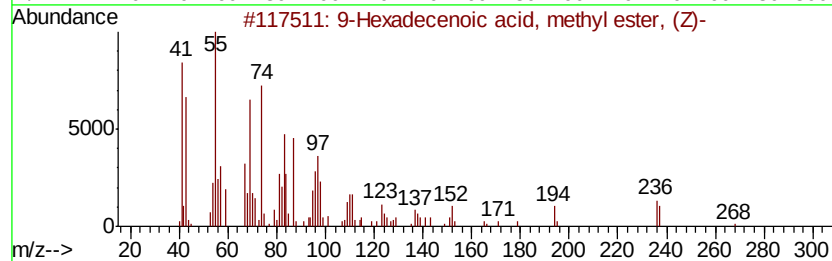
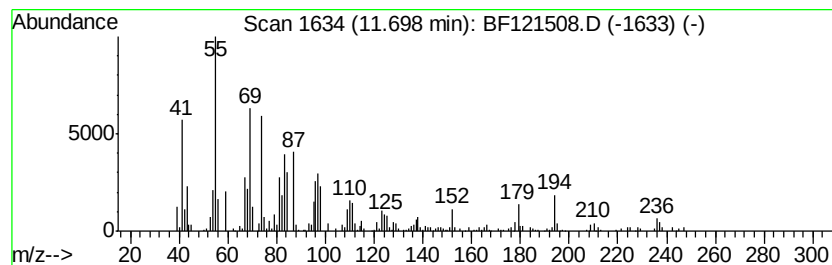
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ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 9-Hexadecenoic acid, methyl... Concentration Rank 8

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 11.70     | 6.05 ng                             | 615597 | Phenanthrene-d10 | 11.39        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 9-Hexadecenoic acid, methyl este... | 268    | C17H32O2         | 001120-25-8  | 99   |
| 2         | 9-Octadecenoic acid (Z)-, methyl... | 296    | C19H36O2         | 000112-62-9  | 93   |
| 3         | 14-Methylpentadec-9-enoic acid m... | 268    | C17H32O2         | 1000365-89-7 | 91   |
| 4         | Methyl hexadec-9-enoate             | 268    | C17H32O2         | 010030-74-7  | 91   |
| 5         | Cyclopropaneoctanoic acid, 2-hex... | 282    | C18H34O2         | 010152-61-1  | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

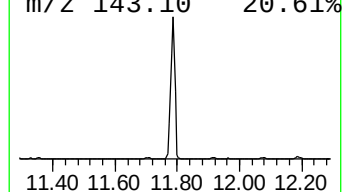
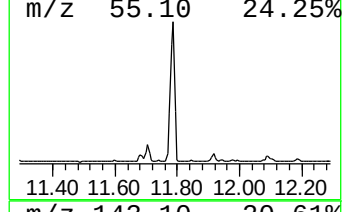
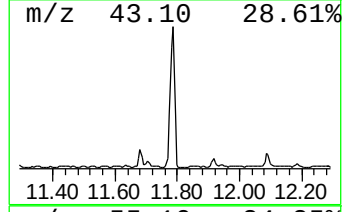
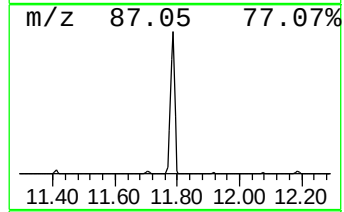
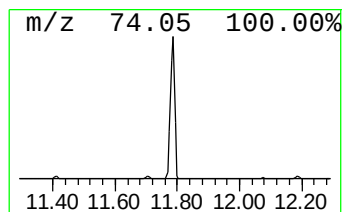
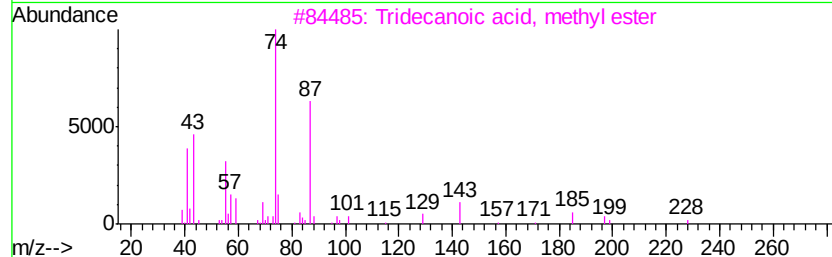
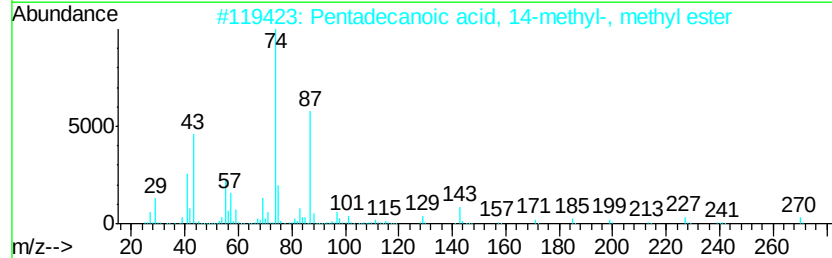
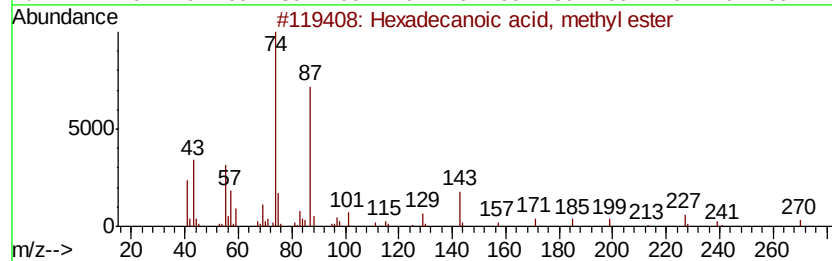
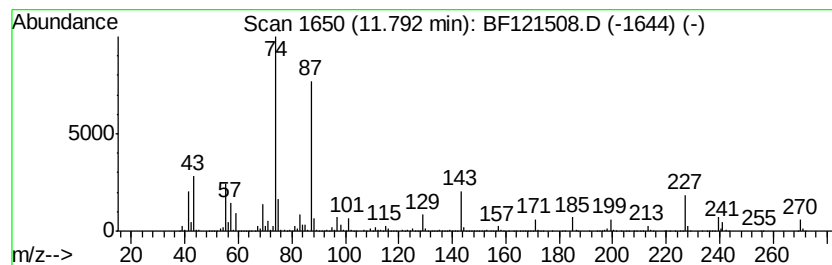
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Hexadecanoic acid, methyl e... Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.79 | 74.21 ng | 7549760 | Phenanthrene-d10 | 11.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 98   |
| 2    |    | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 95   |
| 3    |    | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 90   |
| 4    |    | Pentadecanoic acid, methyl ester    | 256 | C16H32O2 | 007132-64-1 | 90   |
| 5    |    | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
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Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

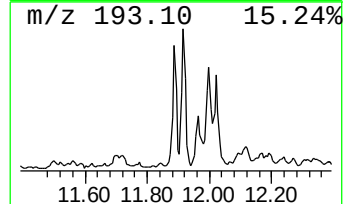
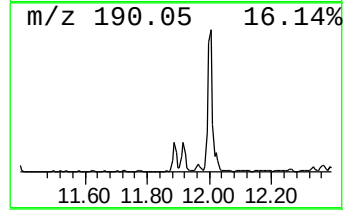
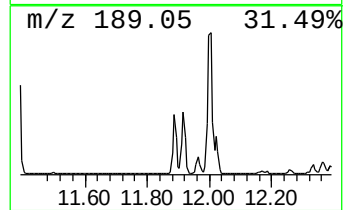
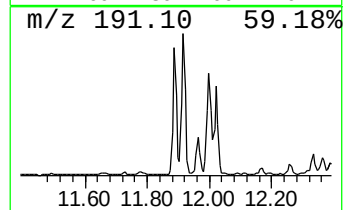
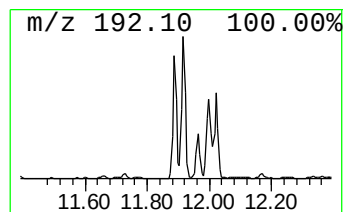
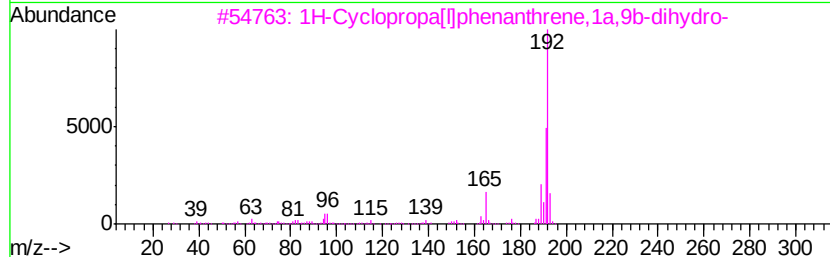
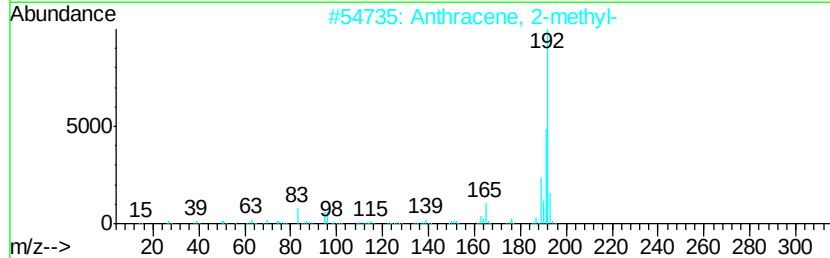
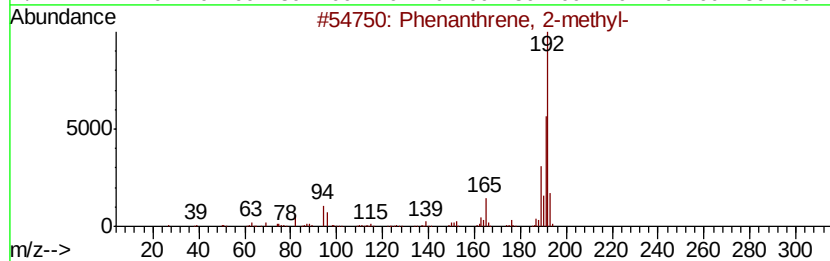
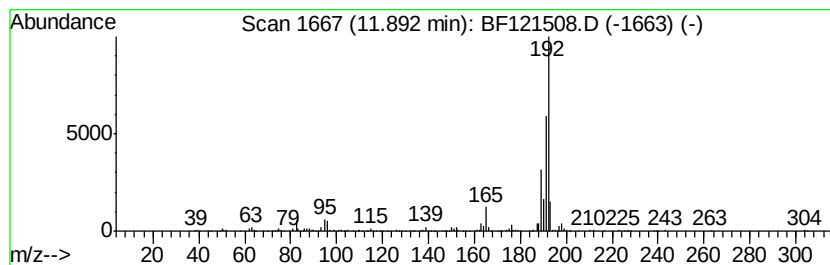
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ClientSampleId :  
OR-02-083120

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 14

| R.T.      | EstConc                               | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------------------------|--------|------------------|-------------|------|
| 11.89     | 3.91 ng                               | 397842 | Phenanthrene-d10 | 11.39       |      |
| Hit# of 5 | Tentative ID                          | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2-methyl-               | 192    | C15H12           | 002531-84-2 | 98   |
| 2         | Anthracene, 2-methyl-                 | 192    | C15H12           | 000613-12-7 | 96   |
| 3         | 1H-Cyclopropa[1,1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 95   |
| 4         | Phenanthrene, 1-methyl-               | 192    | C15H12           | 000832-69-9 | 94   |
| 5         | Anthracene, 1-methyl-                 | 192    | C15H12           | 000610-48-0 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
OR-02-083120

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

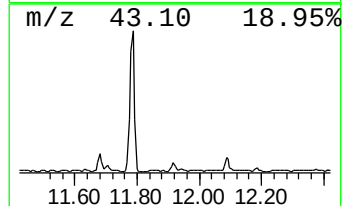
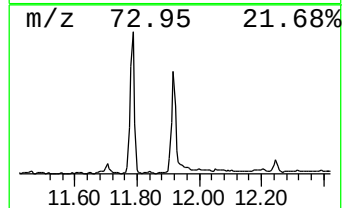
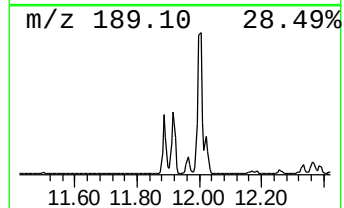
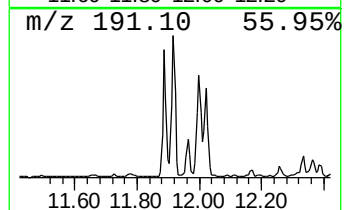
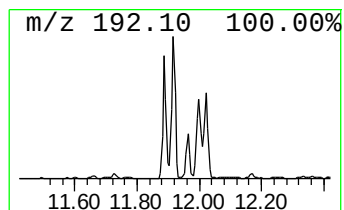
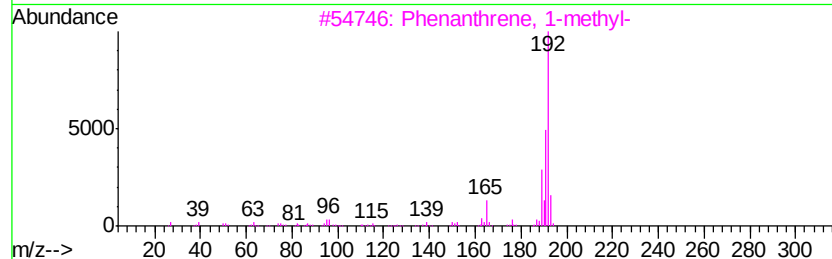
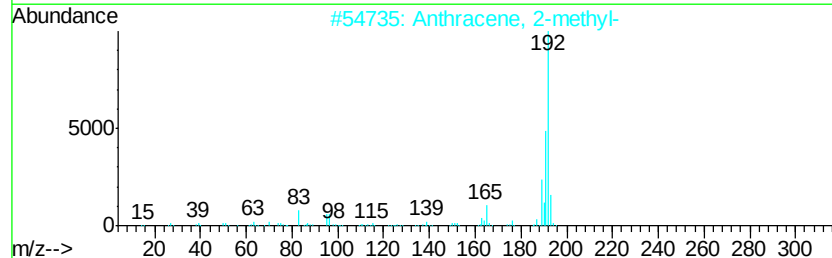
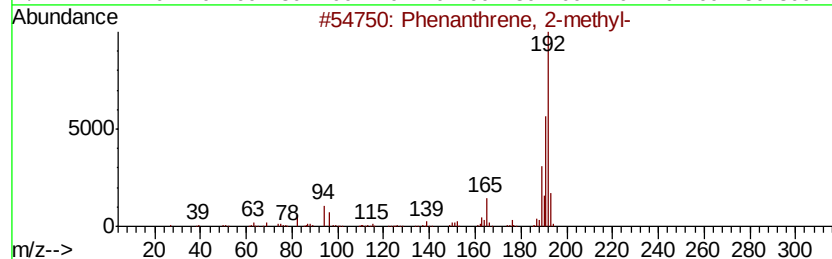
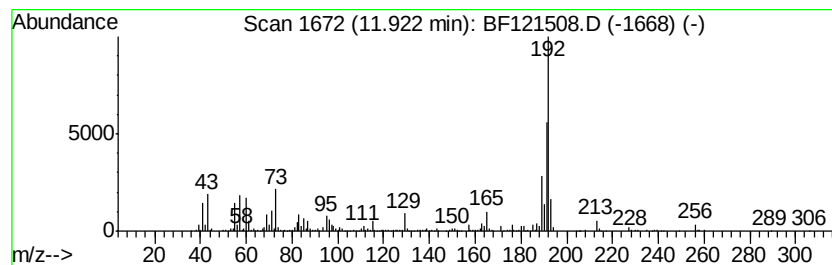
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 Anthracene, 2-methyl- Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.92 | 7.44 ng | 756952 | Phenanthrene-d10 | 11.39 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2-methyl-              | 192 | C15H12  | 002531-84-2 | 97   |
| 2    |    | Anthracene, 2-methyl-                | 192 | C15H12  | 000613-12-7 | 97   |
| 3    |    | Phenanthrene, 1-methyl-              | 192 | C15H12  | 000832-69-9 | 96   |
| 4    |    | Anthracene, 1-methyl-                | 192 | C15H12  | 000610-48-0 | 94   |
| 5    |    | 1H-Cyclopropa[l]phenanthrene, 1a,... | 192 | C15H12  | 000949-41-7 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

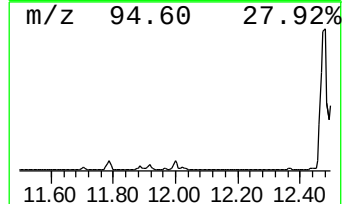
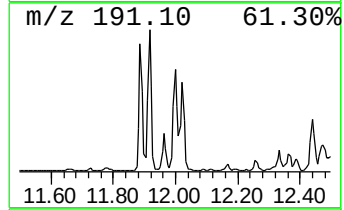
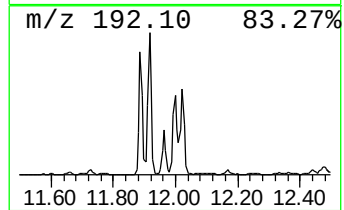
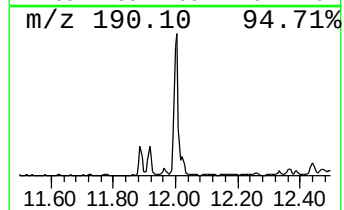
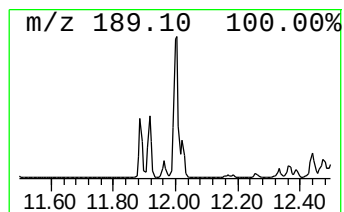
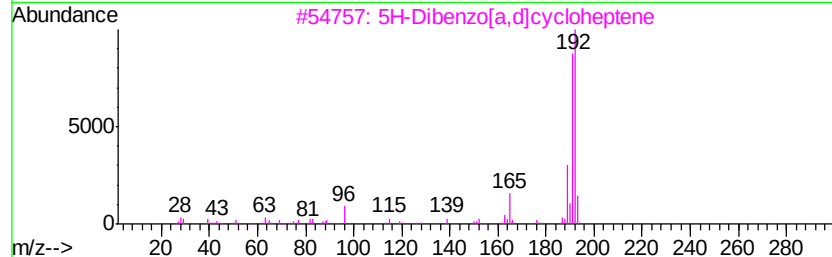
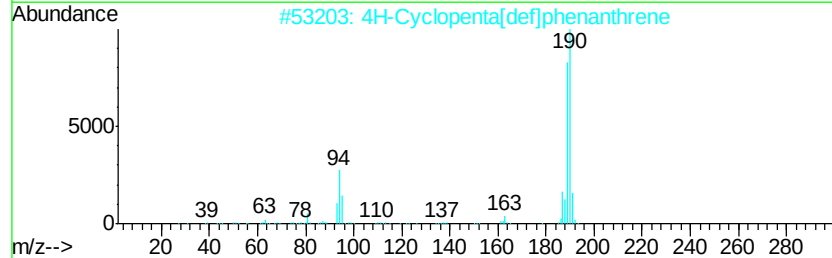
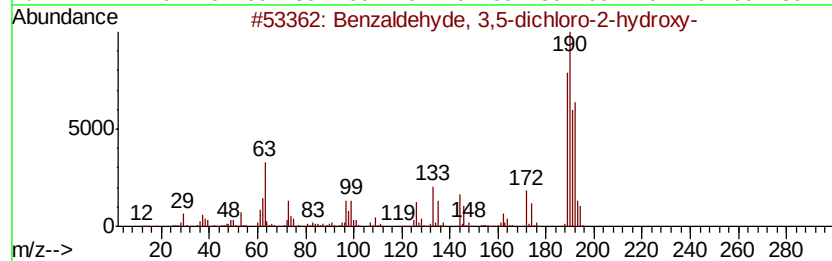
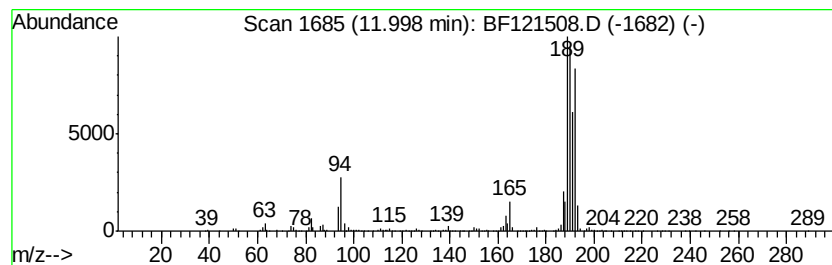
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ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzaldehyde, 3,5-dichloro-... Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 12.00     | 6.69 ng                             | 680143 | Phenanthrene-d10 | 11.39       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Benzaldehyde, 3,5-dichloro-2-hyd... | 190    | C7H4Cl2O2        | 000090-60-8 | 64   |
| 2         | 4H-Cyclopenta[def]phenanthrene      | 190    | C15H10           | 000203-64-5 | 60   |
| 3         | 5H-Dibenzo[a,d]cycloheptene         | 192    | C15H12           | 000256-81-5 | 42   |
| 4         | Naphtho[2,3-b]norborene             | 192    | C15H12           | 107426-38-0 | 38   |
| 5         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

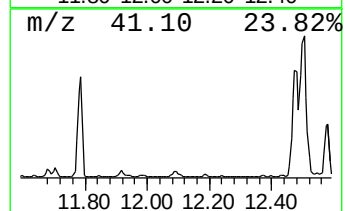
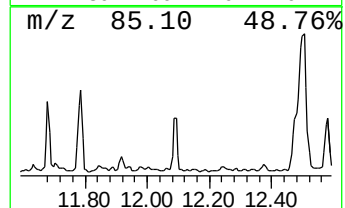
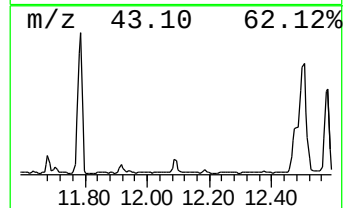
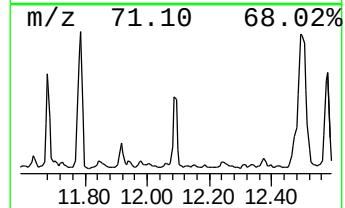
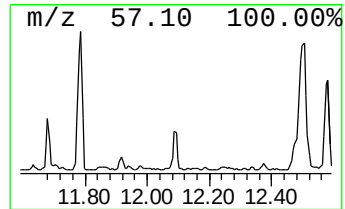
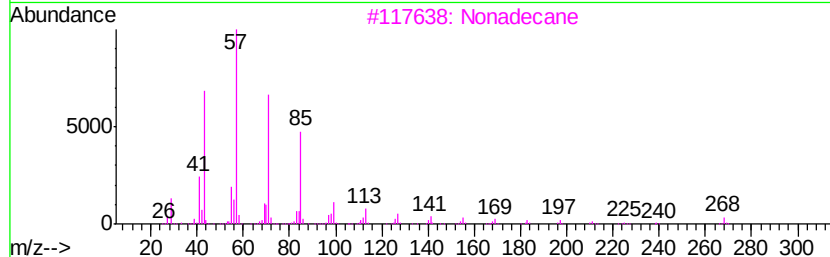
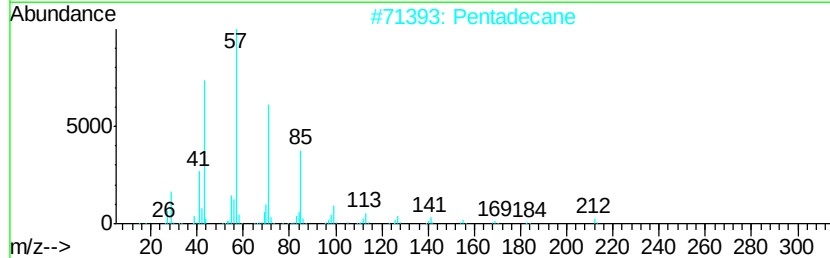
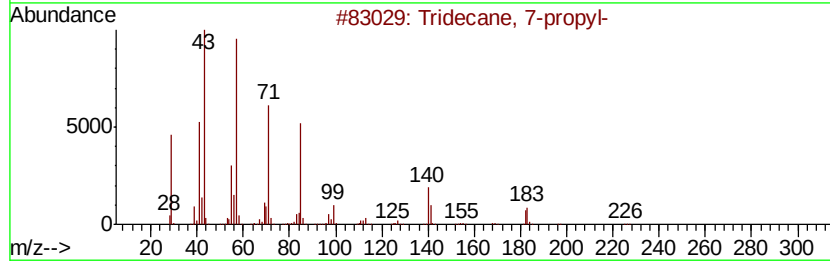
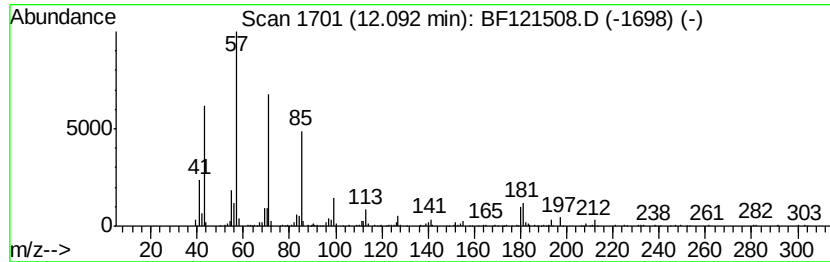
Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Tridecane, 7-propyl- Concentration Rank 10

| R.T.      | EstConc              | Area   | Relative to ISTD | R.T.        |      |
|-----------|----------------------|--------|------------------|-------------|------|
| 12.09     | 5.07 ng              | 515710 | Phenanthrene-d10 | 11.39       |      |
| Hit# of 5 | Tentative ID         | MW     | MolForm          | CAS#        | Qual |
| 1         | Tridecane, 7-propyl- | 226    | C16H34           | 055045-09-5 | 89   |
| 2         | Pentadecane          | 212    | C15H32           | 000629-62-9 | 78   |
| 3         | Nonadecane           | 268    | C19H40           | 000629-92-5 | 68   |
| 4         | Tridecane, 6-propyl- | 226    | C16H34           | 055045-10-8 | 68   |
| 5         | Hexadecane           | 226    | C16H34           | 000544-76-3 | 68   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

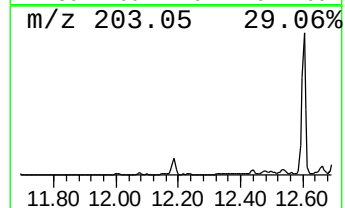
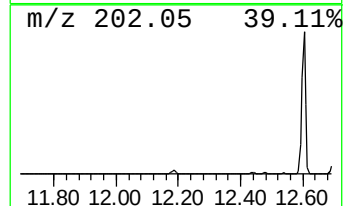
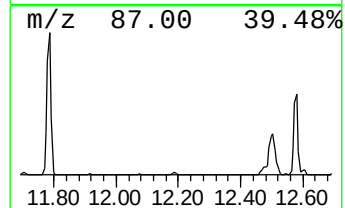
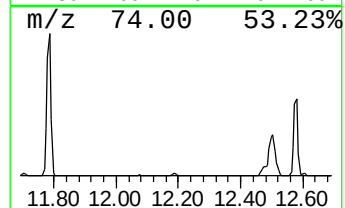
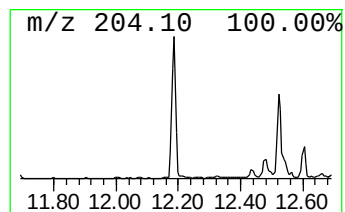
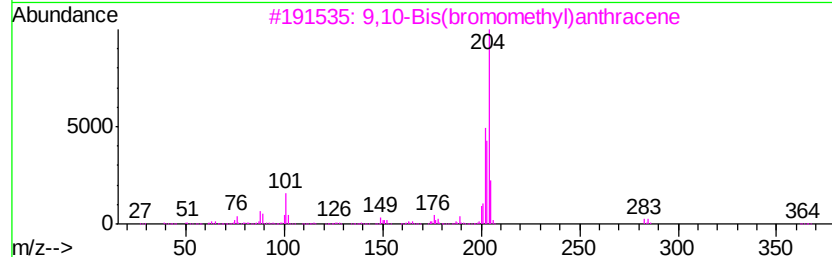
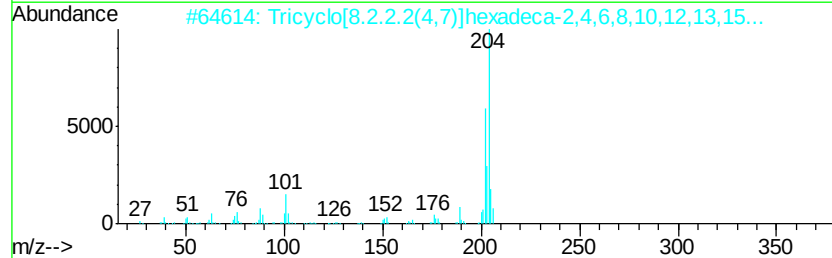
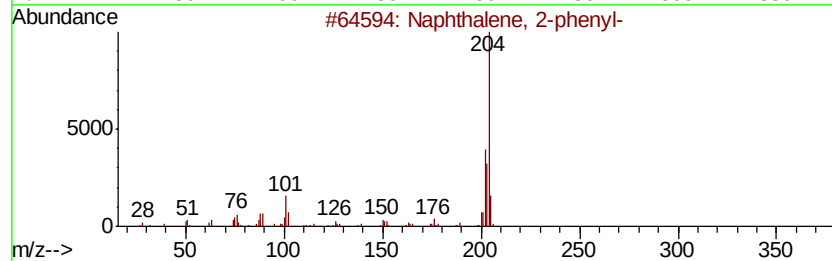
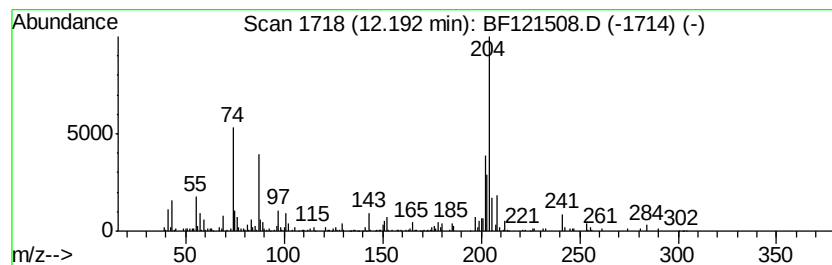
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ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Naphthalene, 2-phenyl- Concentration Rank 15

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 12.19     | 3.91 ng                             | 397311 | Phenanthrene-d10 | 11.39       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2-phenyl-              | 204    | C16H12           | 000612-94-2 | 70   |
| 2         | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204    | C16H12           | 006572-60-7 | 49   |
| 3         | 9,10-Bis(bromomethyl)anthracene     | 362    | C16H12Br2        | 034373-96-1 | 46   |
| 4         | Naphthalene, 1-phenyl-              | 204    | C16H12           | 000605-02-7 | 38   |
| 5         | [1,1'-Biphenyl]-4,4'-dicarbonitrile | 204    | C14H8N2          | 001591-30-6 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

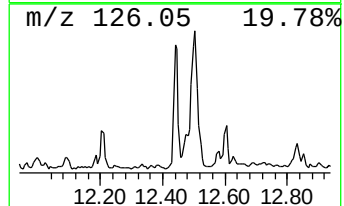
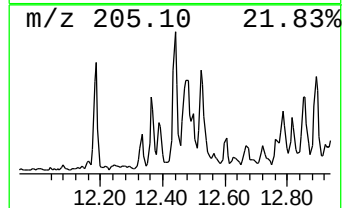
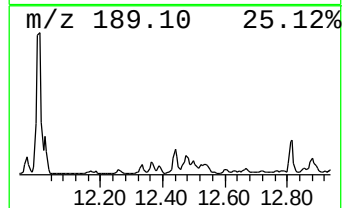
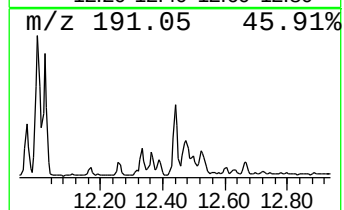
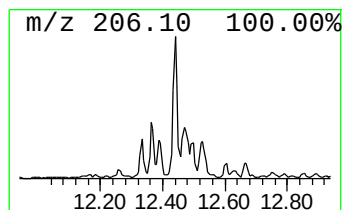
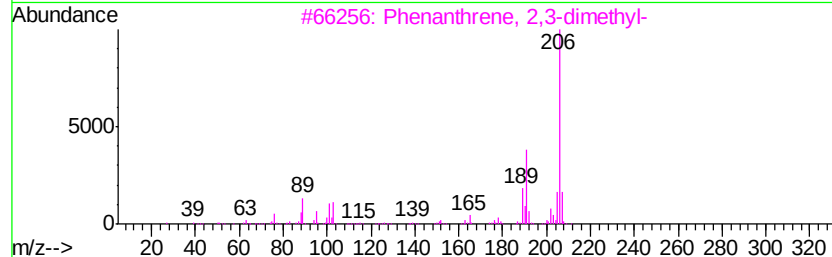
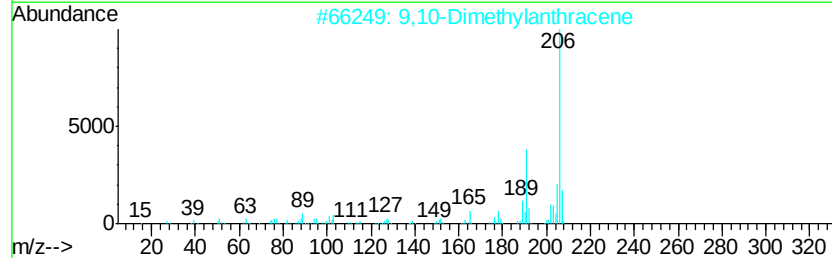
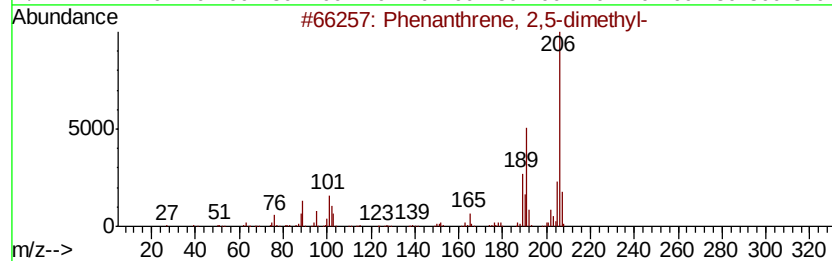
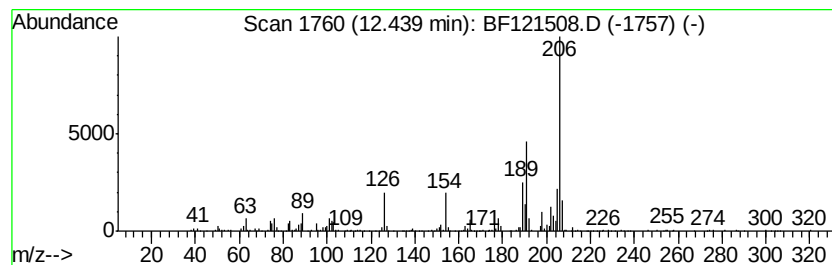
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OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Phenanthrene, 2,5-dimethyl- Concentration Rank 16

| R.T.    | EstConc | Area                        | Relative to ISTD |         | R.T.        |      |
|---------|---------|-----------------------------|------------------|---------|-------------|------|
| 12.44   | 3.72 ng | 378478                      | Phenanthrene-d10 |         | 11.39       |      |
| Hit# of | 5       | Tentative ID                | MW               | MolForm | CAS#        | Qual |
| 1       |         | Phenanthrene, 2,5-dimethyl- | 206              | C16H14  | 003674-66-6 | 94   |
| 2       |         | 9,10-Dimethylantracene      | 206              | C16H14  | 000781-43-1 | 91   |
| 3       |         | Phenanthrene, 2,3-dimethyl- | 206              | C16H14  | 003674-65-5 | 89   |
| 4       |         | Phenanthrene, 1,7-dimethyl- | 206              | C16H14  | 000483-87-4 | 70   |
| 5       |         | Anthracene, 1,4-dimethyl-   | 206              | C16H14  | 000781-92-0 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

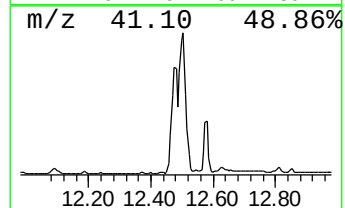
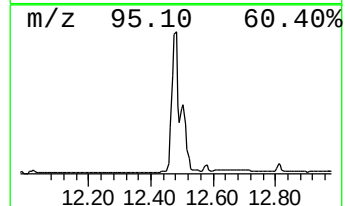
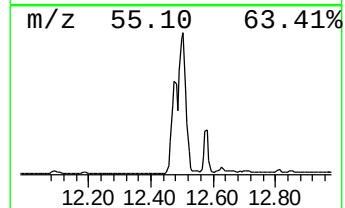
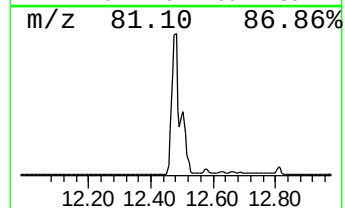
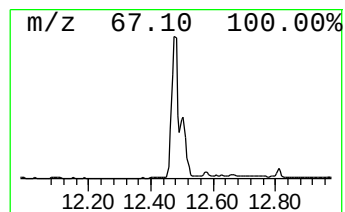
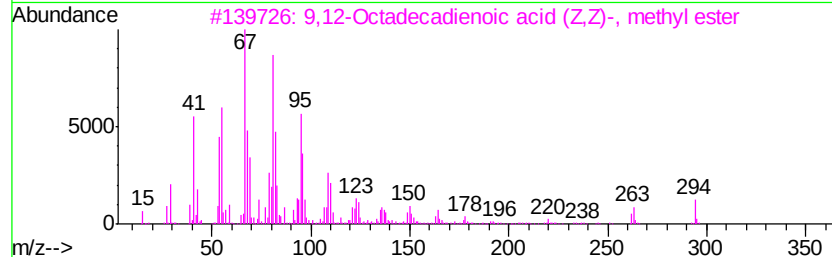
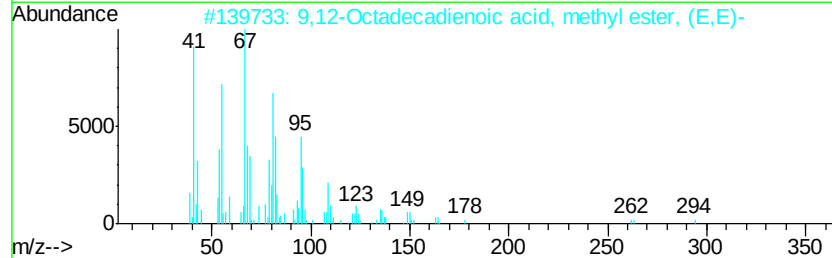
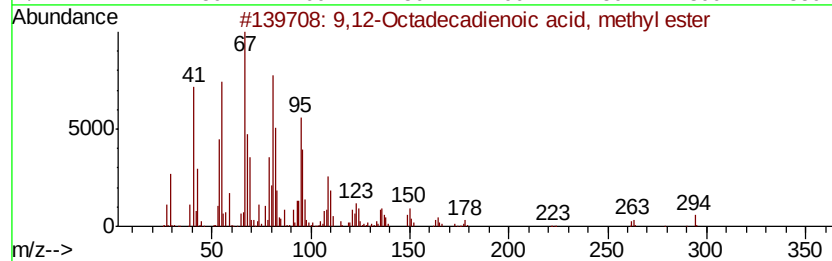
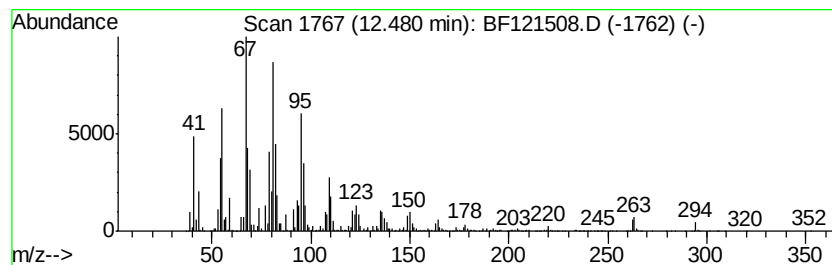
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 9,12-Octadecadienoic acid, ... Concentration Rank 2

| R.T.  | EstConc   | Area     | Relative to ISTD | R.T.  |
|-------|-----------|----------|------------------|-------|
| 12.48 | 117.75 ng | 11979300 | Phenanthrene-d10 | 11.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002462-85-3  | 99   |
| 2    |    | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002566-97-4  | 99   |
| 3    |    | 9,12-Octadecadienoic acid (Z,Z)-... | 294 | C19H34O2 | 000112-63-0  | 99   |
| 4    |    | Methyl 9-cis,11-trans-octadecadi... | 294 | C19H34O2 | 1000336-44-0 | 99   |
| 5    |    | 10,13-Octadecadienoic acid, meth... | 294 | C19H34O2 | 056554-62-2  | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
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Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

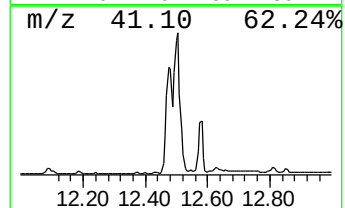
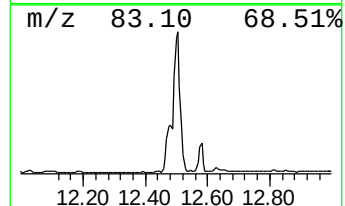
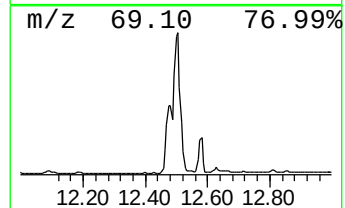
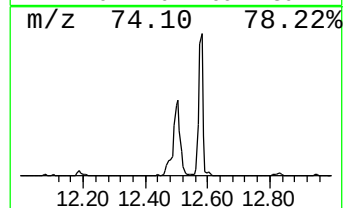
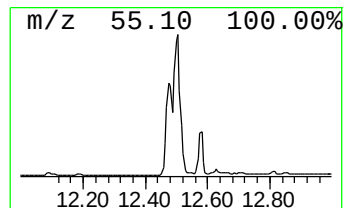
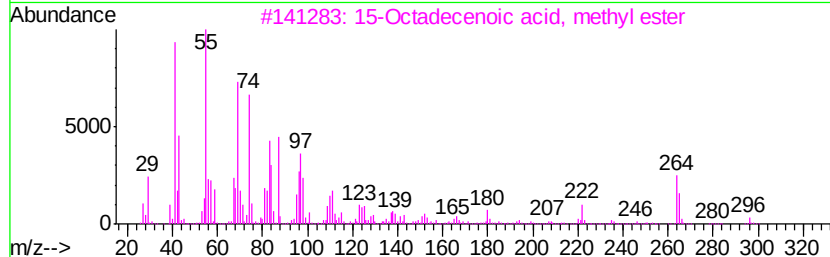
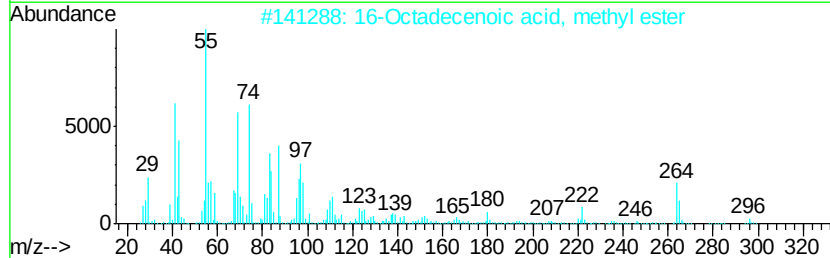
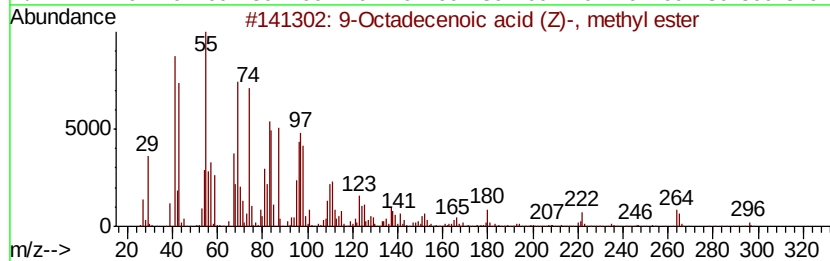
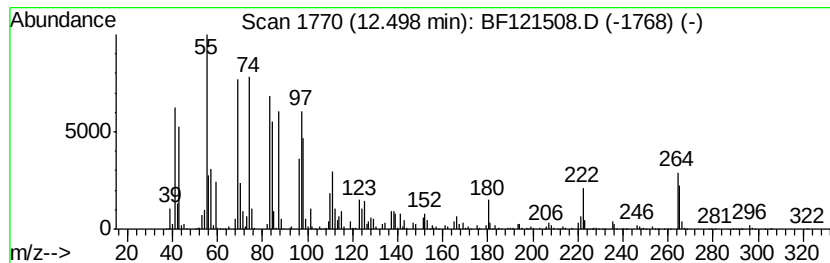
Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 9-Octadecenoic acid (Z)-, m... Concentration Rank 1

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.        |      |      |
|---------|-------------------------------------|--------------|------------------|-------------|------|------|
| 12.50   | 151.00 ng                           | 15362500     | Phenanthrene-d10 | 11.39       |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm     | CAS# | Qual |
| 1       | 9-Octadecenoic acid (Z)-, methyl... | 296          | C19H36O2         | 000112-62-9 | 99   |      |
| 2       | 16-Octadecenoic acid, methyl ester  | 296          | C19H36O2         | 056554-49-5 | 99   |      |
| 3       | 15-Octadecenoic acid, methyl ester  | 296          | C19H36O2         | 004764-72-1 | 96   |      |
| 4       | 9-Octadecenoic acid, methyl este... | 296          | C19H36O2         | 001937-62-8 | 95   |      |
| 5       | 6-Octadecenoic acid, methyl ester   | 296          | C19H36O2         | 052355-31-4 | 95   |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

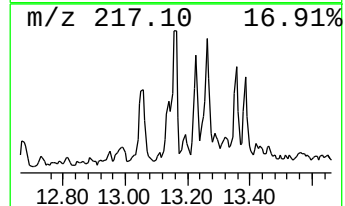
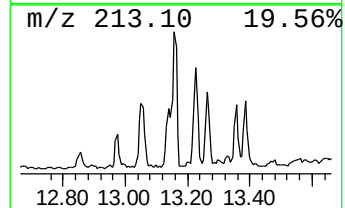
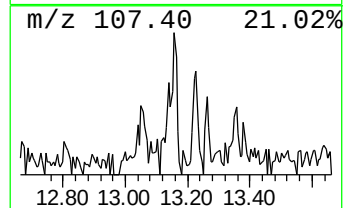
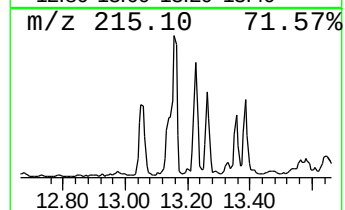
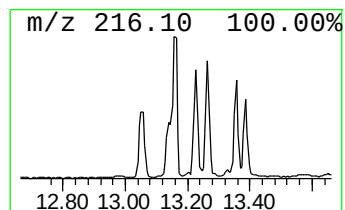
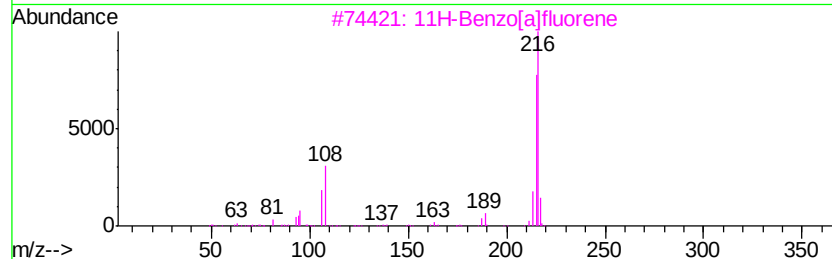
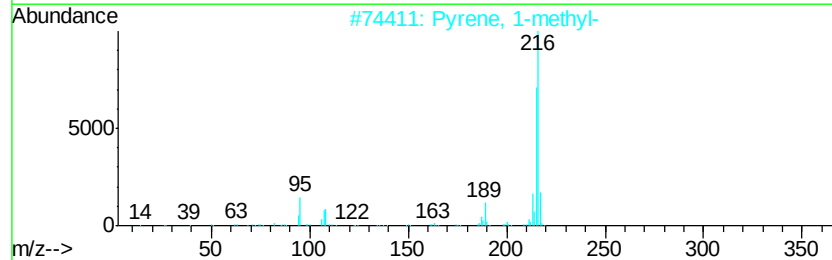
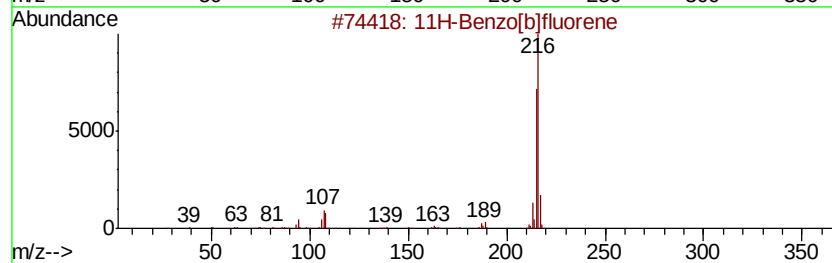
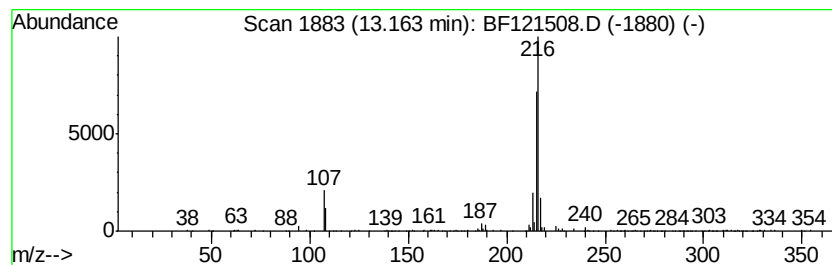
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 11H-Benzo[b]fluorene Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.16 | 2.89 ng | 453802 | Chrysene-d12     | 14.03 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 93   |
| 2    |    | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 87   |
| 3    |    | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 81   |
| 4    |    | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 53   |
| 5    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

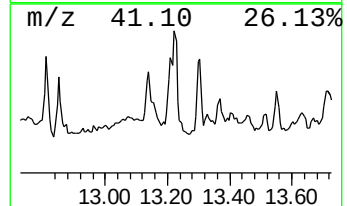
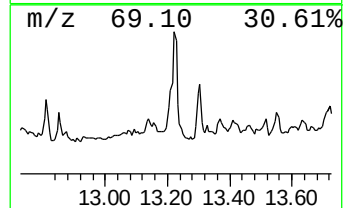
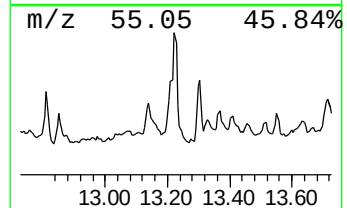
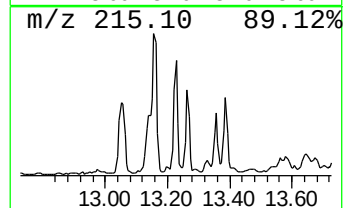
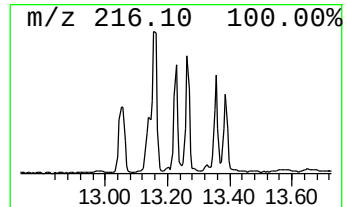
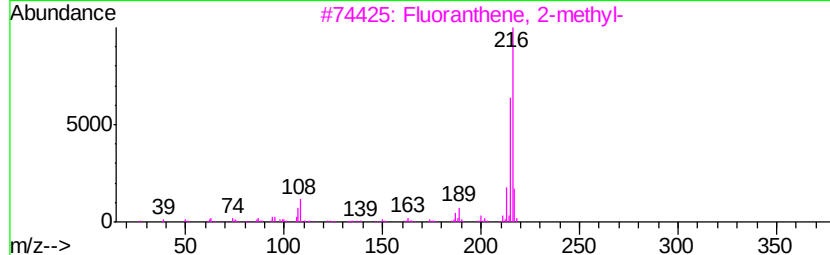
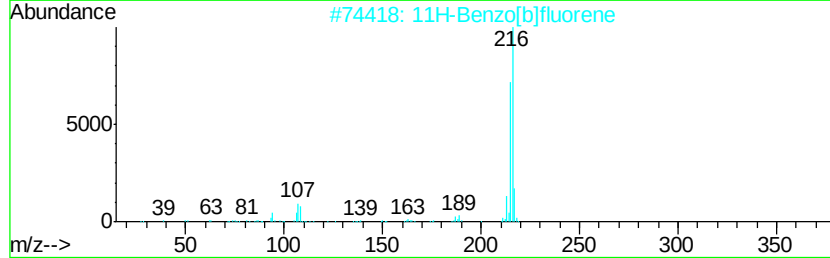
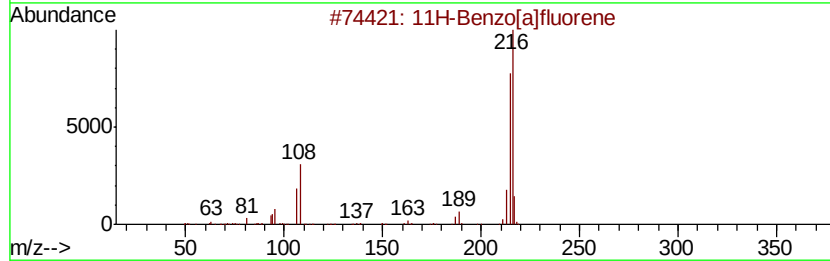
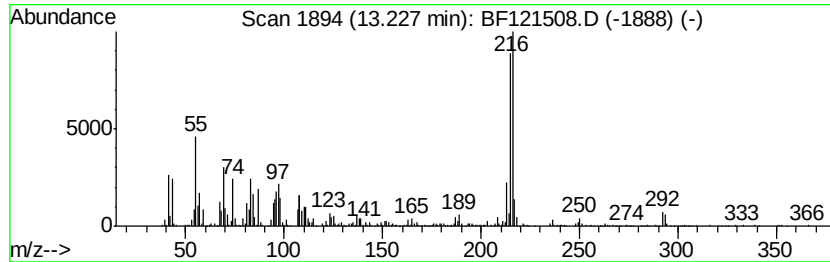
Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 11H-Benzo[a]fluorene Concentration Rank 9

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.           |
|---------|---------|-------------------------|------------------|----------------|
| 13.23   | 6.01 ng | 942708                  | Chrysene-d12     | 14.03          |
| Hit# of | 5       | Tentative ID            | MW MolForm       | CAS# Qual      |
| 1       |         | 11H-Benzo[a]fluorene    | 216 C17H12       | 000238-84-6 62 |
| 2       |         | 11H-Benzo[b]fluorene    | 216 C17H12       | 000243-17-4 55 |
| 3       |         | Fluoranthene, 2-methyl- | 216 C17H12       | 033543-31-6 50 |
| 4       |         | Pyrene, 1-methyl-       | 216 C17H12       | 002381-21-7 49 |
| 5       |         | Pyrene, 2-methyl-       | 216 C17H12       | 003442-78-2 43 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

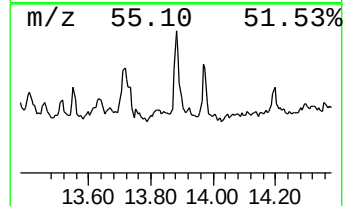
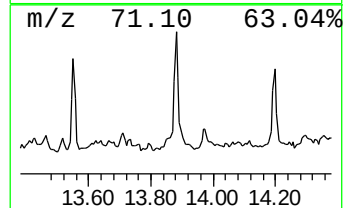
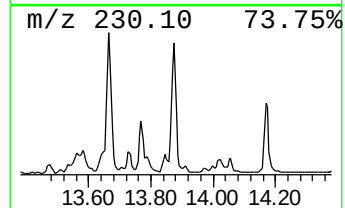
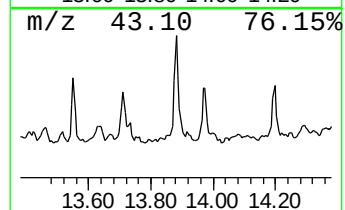
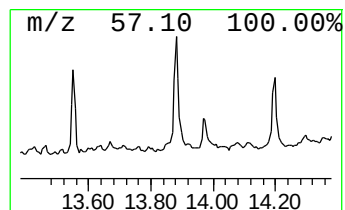
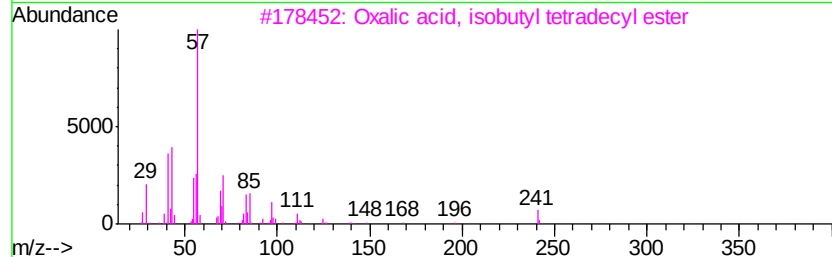
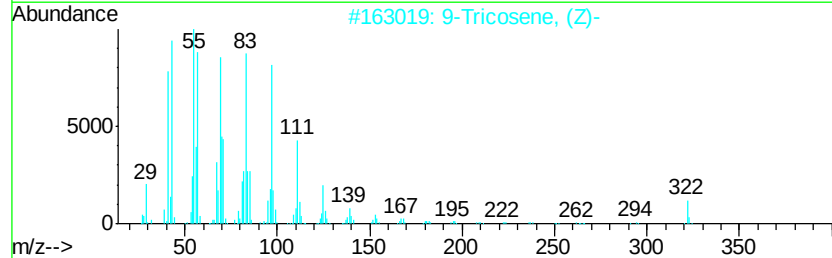
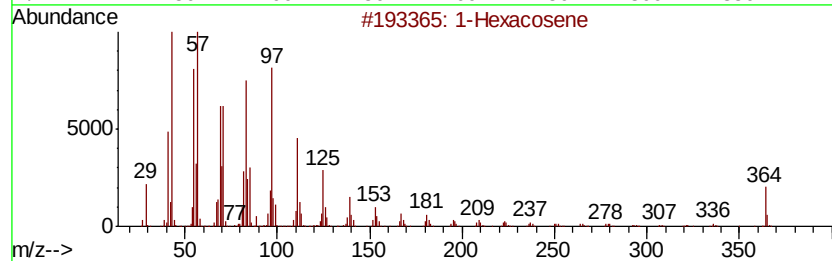
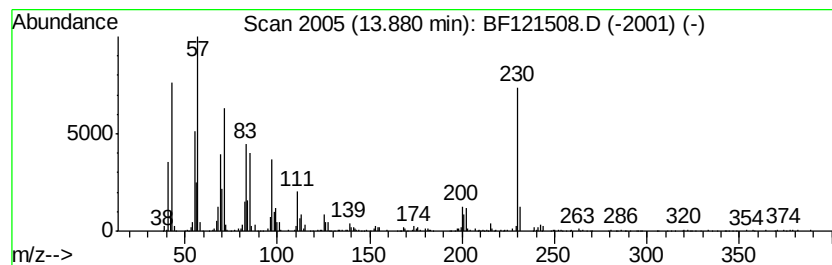
Instrument :  
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ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 1-Hexacosene Concentration Rank 17

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 13.88     | 3.29 ng                             | 516187 | Chrysene-d12     | 14.03        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-Hexacosene                        | 364    | C26H52           | 018835-33-1  | 93   |
| 2         | 9-Tricosene, (Z)-                   | 322    | C23H46           | 027519-02-4  | 92   |
| 3         | Oxalic acid, isobutyl tetradecyl... | 342    | C20H38O4         | 1000309-37-9 | 60   |
| 4         | 1-Decanol, 2-hexyl-                 | 242    | C16H34O          | 002425-77-6  | 60   |
| 5         | Oxalic acid, isobutyl hexadecyl ... | 370    | C22H42O4         | 1000309-38-1 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\  
Data File : BF121508.D  
Acq On : 1 Sep 2020 21:13  
Operator : JU/CG  
Sample : L3506-11 5X  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

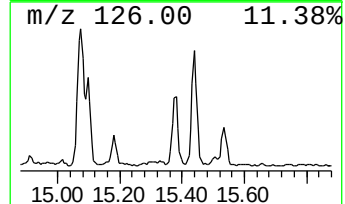
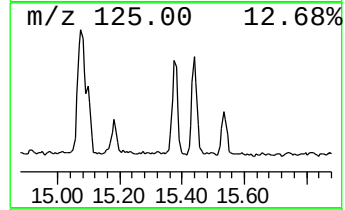
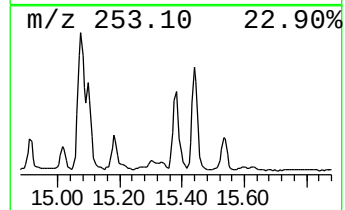
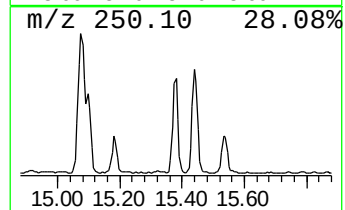
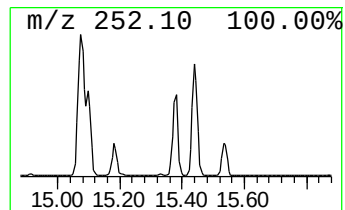
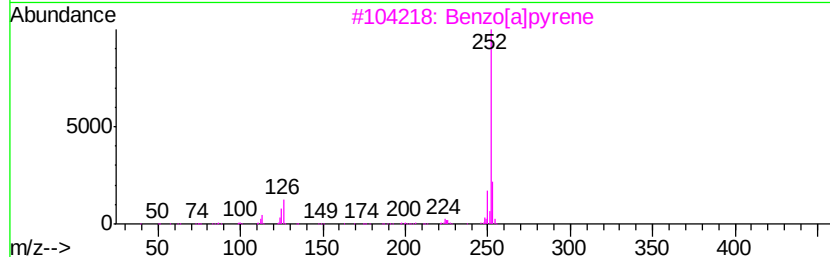
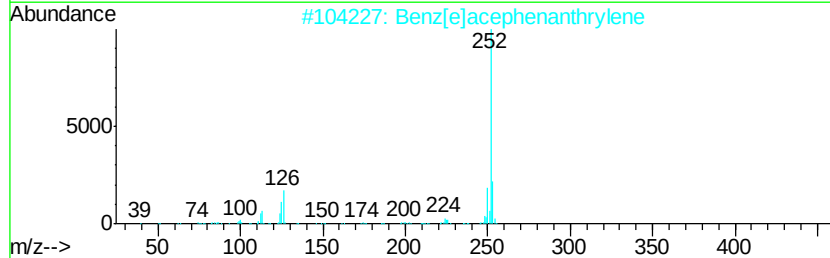
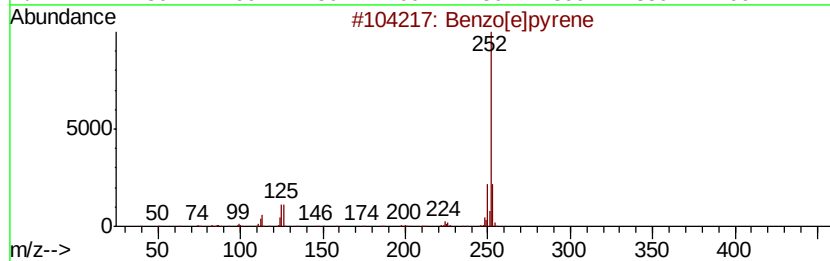
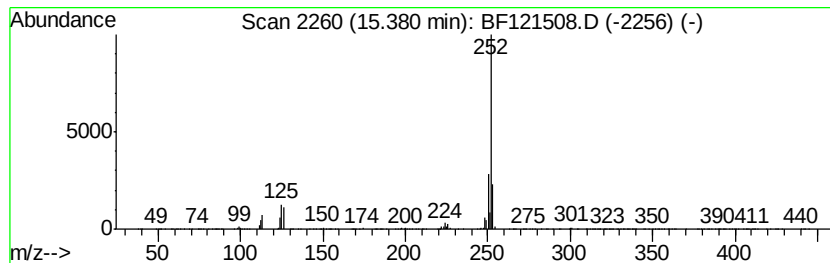
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Benzo[e]pyrene Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.38 | 7.28 ng | 674694 | Perylene-d12     | 15.51 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[e]pyrene            | 252 | C20H12  | 000192-97-2 | 98   |
| 2    |    | Benzo[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 93   |
| 3    |    | Benzo[a]pyrene            | 252 | C20H12  | 000050-32-8 | 93   |
| 4    |    | Benzo[k]fluoranthene      | 252 | C20H12  | 000207-08-9 | 93   |
| 5    |    | Benzo[j]fluoranthene      | 252 | C20H12  | 000205-82-3 | 90   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF090120\  
 Data File : BF121508.D  
 Acq On : 1 Sep 2020 21:13  
 Operator : JU/CG  
 Sample : L3506-11 5X  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OR-02-083120

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF082720.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Naphthalene, 2,3-... | 9.56  | 3.2     | ng    | 365603   | 3                     | 9.90  | 2314600 | 20.0 |
| Hexadecane           | 10.33 | 2.8     | ng    | 323414   | 3                     | 9.90  | 2314600 | 20.0 |
| Heptadecane          | 10.80 | 6.5     | ng    | 657548   | 4                     | 11.39 | 2034750 | 20.0 |
| Bacchotricuneatin c  | 11.25 | 4.3     | ng    | 436202   | 4                     | 11.39 | 2034750 | 20.0 |
| Naphtho[1,2-b]thi... | 11.28 | 4.3     | ng    | 432306   | 4                     | 11.39 | 2034750 | 20.0 |
| Tetracosane          | 11.68 | 4.1     | ng    | 421697   | 4                     | 11.39 | 2034750 | 20.0 |
| 9-Hexadecenoic ac... | 11.70 | 6.0     | ng    | 615597   | 4                     | 11.39 | 2034750 | 20.0 |
| Hexadecanoic acid... | 11.79 | 74.2    | ng    | 7549760  | 4                     | 11.39 | 2034750 | 20.0 |
| Phenanthrene, 2-m... | 11.89 | 3.9     | ng    | 397842   | 4                     | 11.39 | 2034750 | 20.0 |
| Anthracene, 2-met... | 11.92 | 7.4     | ng    | 756952   | 4                     | 11.39 | 2034750 | 20.0 |
| Benzaldehyde, 3,5... | 12.00 | 6.7     | ng    | 680143   | 4                     | 11.39 | 2034750 | 20.0 |
| Tridecane, 7-propyl- | 12.09 | 5.1     | ng    | 515710   | 4                     | 11.39 | 2034750 | 20.0 |
| Naphthalene, 2-ph... | 12.19 | 3.9     | ng    | 397311   | 4                     | 11.39 | 2034750 | 20.0 |
| Phenanthrene, 2,5... | 12.44 | 3.7     | ng    | 378478   | 4                     | 11.39 | 2034750 | 20.0 |
| 9,12-Octadecadien... | 12.48 | 117.8   | ng    | 11979300 | 4                     | 11.39 | 2034750 | 20.0 |
| 9-Octadecenoic ac... | 12.50 | 151.0   | ng    | 15362500 | 4                     | 11.39 | 2034750 | 20.0 |
| 11H-Benzo[b]fluorene | 13.16 | 2.9     | ng    | 453802   | 5                     | 14.03 | 3139480 | 20.0 |
| 11H-Benzo[a]fluorene | 13.23 | 6.0     | ng    | 942708   | 5                     | 14.03 | 3139480 | 20.0 |
| 1-Hexacosene         | 13.88 | 3.3     | ng    | 516187   | 5                     | 14.03 | 3139480 | 20.0 |
| Benzo[e]pyrene       | 15.38 | 7.3     | ng    | 674694   | 6                     | 15.51 | 1853570 | 20.0 |