

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF083120\  
 Data File : BF121486.D  
 Acq On : 31 Aug 2020 19:32  
 Operator : JU/CG  
 Sample : L3847-15  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 LINDEN-JOP-BH-10-B

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         | 2.146       | 5             | 10          | 23           | rVB      | 1109441        | 1431185       | 26.69%          | 1.679%        |
| 2         | 5.069       | 504           | 507         | 521          | rVB      | 52712          | 77516         | 1.45%           | 0.091%        |
| 3         | 5.475       | 571           | 576         | 579          | rBV      | 3396303        | 3075060       | 57.35%          | 3.607%        |
| 4         | 6.481       | 743           | 747         | 750          | rBV      | 3536566        | 3161736       | 58.97%          | 3.708%        |
| 5         | 6.863       | 808           | 812         | 815          | rBV      | 2154445        | 1842651       | 34.37%          | 2.161%        |
| 6         | 7.422       | 902           | 907         | 910          | rBV      | 2464589        | 2175251       | 40.57%          | 2.551%        |
| 7         | 8.145       | 1026          | 1030        | 1032         | rBV      | 2762782        | 2294949       | 42.80%          | 2.692%        |
| 8         | 8.163       | 1032          | 1033        | 1041         | rVB      | 315017         | 237427        | 4.43%           | 0.278%        |
| 9         | 8.857       | 1148          | 1151        | 1157         | rVB      | 192070         | 151102        | 2.82%           | 0.177%        |
| 10        | 8.957       | 1163          | 1168        | 1173         | rBV      | 163913         | 139450        | 2.60%           | 0.164%        |
| 11        | 9.222       | 1208          | 1213        | 1216         | rBV      | 4769992        | 4144421       | 77.30%          | 4.861%        |
| 12        | 9.322       | 1227          | 1230        | 1235         | rVB      | 77436          | 87348         | 1.63%           | 0.102%        |
| 13        | 9.486       | 1252          | 1258        | 1262         | rBV      | 82874          | 124798        | 2.33%           | 0.146%        |
| 14        | 9.557       | 1265          | 1270        | 1272         | rVV      | 163747         | 154726        | 2.89%           | 0.181%        |
| 15        | 9.586       | 1272          | 1275        | 1277         | rVV      | 88321          | 87165         | 1.63%           | 0.102%        |
| 16        | 9.610       | 1277          | 1279        | 1284         | rVV      | 194482         | 185472        | 3.46%           | 0.218%        |
| 17        | 9.675       | 1286          | 1290        | 1296         | rVV      | 86050          | 97897         | 1.83%           | 0.115%        |
| 18        | 9.757       | 1299          | 1304        | 1308         | rBV      | 236889         | 211204        | 3.94%           | 0.248%        |
| 19        | 9.898       | 1322          | 1328        | 1331         | rBV      | 2865944        | 2534019       | 47.26%          | 2.972%        |
| 20        | 9.933       | 1331          | 1334        | 1337         | rVB      | 571525         | 491609        | 9.17%           | 0.577%        |
| 21        | 10.057      | 1350          | 1355        | 1359         | rBV2     | 54713          | 67399         | 1.26%           | 0.079%        |
| 22        | 10.104      | 1359          | 1363        | 1367         | rVV      | 464053         | 424023        | 7.91%           | 0.497%        |
| 23        | 10.216      | 1378          | 1382        | 1385         | rBV      | 71049          | 76819         | 1.43%           | 0.090%        |
| 24        | 10.304      | 1392          | 1397        | 1399         | rBV3     | 76441          | 95971         | 1.79%           | 0.113%        |
| 25        | 10.445      | 1417          | 1421        | 1426         | rBV      | 793202         | 666772        | 12.44%          | 0.782%        |
| 26        | 10.604      | 1445          | 1448        | 1454         | rVB      | 200676         | 187013        | 3.49%           | 0.219%        |
| 27        | 10.686      | 1454          | 1462        | 1466         | rBV      | 2466042        | 2361990       | 44.05%          | 2.770%        |
| 28        | 10.733      | 1466          | 1470        | 1474         | rVB2     | 59231          | 82966         | 1.55%           | 0.097%        |
| 29        | 10.810      | 1479          | 1483        | 1489         | rVB4     | 93231          | 121527        | 2.27%           | 0.143%        |
| 30        | 10.992      | 1508          | 1514        | 1517         | rBV3     | 174351         | 253652        | 4.73%           | 0.298%        |
| 31        | 11.022      | 1517          | 1519        | 1522         | rVV2     | 98973          | 90051         | 1.68%           | 0.106%        |
| 32        | 11.074      | 1525          | 1528        | 1531         | rVV      | 77986          | 81713         | 1.52%           | 0.096%        |
| 33        | 11.122      | 1531          | 1536        | 1538         | rVV3     | 112247         | 143979        | 2.69%           | 0.169%        |
| 34        | 11.145      | 1538          | 1540        | 1545         | rVV2     | 135353         | 150453        | 2.81%           | 0.176%        |

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 Sample : L3847-15  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 LINDEN-JOP-BH-10-B

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 | 11.186 | 1545 | 1547 | 1551 | rVV  | 177698  | 164832  | 3.07%   | 0.193% |
| 36 | 11.233 | 1551 | 1555 | 1560 | rVV4 | 124381  | 234546  | 4.37%   | 0.275% |
| 37 | 11.280 | 1560 | 1563 | 1569 | rVB  | 389889  | 347729  | 6.49%   | 0.408% |
| 38 | 11.386 | 1577 | 1581 | 1583 | rBV  | 2571909 | 2473382 | 46.13%  | 2.901% |
| 39 | 11.410 | 1583 | 1585 | 1588 | rVV  | 5108431 | 4545155 | 84.77%  | 5.331% |
| 40 | 11.463 | 1588 | 1594 | 1597 | rVV  | 1570606 | 1438526 | 26.83%  | 1.687% |
| 41 | 11.492 | 1597 | 1599 | 1602 | rVB  | 168618  | 147464  | 2.75%   | 0.173% |
| 42 | 11.610 | 1614 | 1619 | 1622 | rBV2 | 536939  | 608108  | 11.34%  | 0.713% |
| 43 | 11.680 | 1629 | 1631 | 1634 | rBV2 | 145450  | 110524  | 2.06%   | 0.130% |
| 44 | 11.716 | 1634 | 1637 | 1641 | rVB3 | 110614  | 93122   | 1.74%   | 0.109% |
| 45 | 11.798 | 1648 | 1651 | 1656 | rVB  | 67107   | 79278   | 1.48%   | 0.093% |
| 46 | 11.886 | 1660 | 1666 | 1668 | rBV  | 671305  | 603531  | 11.26%  | 0.708% |
| 47 | 11.916 | 1668 | 1671 | 1675 | rVV  | 1084026 | 1036934 | 19.34%  | 1.216% |
| 48 | 11.963 | 1675 | 1679 | 1681 | rVV  | 388793  | 342659  | 6.39%   | 0.402% |
| 49 | 11.998 | 1681 | 1685 | 1687 | rVV  | 1048317 | 1068198 | 19.92%  | 1.253% |
| 50 | 12.021 | 1687 | 1689 | 1692 | rVB  | 424028  | 336030  | 6.27%   | 0.394% |
| 51 | 12.086 | 1698 | 1700 | 1703 | rVB2 | 99126   | 84634   | 1.58%   | 0.099% |
| 52 | 12.180 | 1713 | 1716 | 1718 | rBV  | 469590  | 385143  | 7.18%   | 0.452% |
| 53 | 12.204 | 1718 | 1720 | 1723 | rVB  | 272493  | 210608  | 3.93%   | 0.247% |
| 54 | 12.333 | 1739 | 1742 | 1744 | rVB  | 132480  | 103062  | 1.92%   | 0.121% |
| 55 | 12.363 | 1744 | 1747 | 1749 | rBV  | 108048  | 90400   | 1.69%   | 0.106% |
| 56 | 12.386 | 1749 | 1751 | 1753 | rVB  | 129996  | 98260   | 1.83%   | 0.115% |
| 57 | 12.433 | 1756 | 1759 | 1762 | rBV  | 312679  | 349563  | 6.52%   | 0.410% |
| 58 | 12.474 | 1763 | 1766 | 1771 | rVB3 | 234220  | 305312  | 5.69%   | 0.358% |
| 59 | 12.521 | 1771 | 1774 | 1779 | rVB2 | 255008  | 288097  | 5.37%   | 0.338% |
| 60 | 12.604 | 1783 | 1788 | 1791 | rBV  | 5531645 | 5087734 | 94.89%  | 5.967% |
| 61 | 12.692 | 1801 | 1803 | 1806 | rBV2 | 144274  | 140273  | 2.62%   | 0.165% |
| 62 | 12.751 | 1810 | 1813 | 1817 | rVB  | 186745  | 170066  | 3.17%   | 0.199% |
| 63 | 12.833 | 1820 | 1827 | 1832 | rBV2 | 4315007 | 5361680 | 100.00% | 6.289% |
| 64 | 12.874 | 1832 | 1834 | 1839 | rVB2 | 247862  | 280386  | 5.23%   | 0.329% |
| 65 | 12.945 | 1843 | 1846 | 1847 | rBV  | 303933  | 249281  | 4.65%   | 0.292% |
| 66 | 12.968 | 1847 | 1850 | 1857 | rVB  | 3624657 | 3709527 | 69.19%  | 4.351% |
| 67 | 13.045 | 1858 | 1863 | 1866 | rBV2 | 332663  | 558462  | 10.42%  | 0.655% |
| 68 | 13.074 | 1866 | 1868 | 1871 | rVB  | 138467  | 98922   | 1.84%   | 0.116% |
| 69 | 13.157 | 1875 | 1882 | 1885 | rVB2 | 724892  | 1002848 | 18.70%  | 1.176% |
| 70 | 13.186 | 1885 | 1887 | 1889 | rBV2 | 73518   | 83411   | 1.56%   | 0.098% |
| 71 | 13.221 | 1890 | 1893 | 1896 | rVB  | 527951  | 430747  | 8.03%   | 0.505% |

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 Sample : L3847-15  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 LINDEN-JOP-BH-10-B

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  | 13.257 | 1896 | 1899 | 1902 | rBV2 | 493787  | 490474  | 9.15%  | 0.575% |
| 73  | 13.286 | 1902 | 1904 | 1907 | rVB  | 100196  | 78031   | 1.46%  | 0.092% |
| 74  | 13.315 | 1907 | 1909 | 1912 | rBV3 | 101208  | 108079  | 2.02%  | 0.127% |
| 75  | 13.351 | 1912 | 1915 | 1918 | rBV  | 288217  | 220551  | 4.11%  | 0.259% |
| 76  | 13.380 | 1918 | 1920 | 1924 | rBV2 | 285301  | 270478  | 5.04%  | 0.317% |
| 77  | 13.457 | 1930 | 1933 | 1938 | rVB4 | 153912  | 178768  | 3.33%  | 0.210% |
| 78  | 13.533 | 1943 | 1946 | 1947 | rBV  | 93688   | 111228  | 2.07%  | 0.130% |
| 79  | 13.663 | 1965 | 1968 | 1972 | rVB  | 290840  | 330503  | 6.16%  | 0.388% |
| 80  | 13.774 | 1983 | 1987 | 1989 | rBV2 | 387796  | 468748  | 8.74%  | 0.550% |
| 81  | 13.804 | 1989 | 1992 | 1994 | rVV  | 497615  | 498115  | 9.29%  | 0.584% |
| 82  | 13.827 | 1994 | 1996 | 2000 | rVV2 | 373773  | 531560  | 9.91%  | 0.623% |
| 83  | 13.868 | 2000 | 2003 | 2010 | rVB3 | 771333  | 879208  | 16.40% | 1.031% |
| 84  | 14.021 | 2022 | 2029 | 2032 | rBV3 | 3277047 | 4825894 | 90.01% | 5.660% |
| 85  | 14.051 | 2032 | 2034 | 2040 | rVB  | 2337403 | 2075922 | 38.72% | 2.435% |
| 86  | 14.133 | 2045 | 2048 | 2050 | rVV  | 335545  | 259795  | 4.85%  | 0.305% |
| 87  | 14.245 | 2064 | 2067 | 2071 | rBV2 | 231836  | 311419  | 5.81%  | 0.365% |
| 88  | 14.410 | 2093 | 2095 | 2098 | rVB  | 395293  | 314842  | 5.87%  | 0.369% |
| 89  | 14.445 | 2099 | 2101 | 2104 | rVV  | 279630  | 226445  | 4.22%  | 0.266% |
| 90  | 14.504 | 2108 | 2111 | 2114 | rVV2 | 309431  | 352489  | 6.57%  | 0.413% |
| 91  | 14.533 | 2114 | 2116 | 2118 | rVV  | 253708  | 248460  | 4.63%  | 0.291% |
| 92  | 14.557 | 2118 | 2120 | 2124 | rVB3 | 276658  | 256537  | 4.78%  | 0.301% |
| 93  | 15.068 | 2203 | 2207 | 2210 | rBV  | 1839417 | 2699730 | 50.35% | 3.166% |
| 94  | 15.174 | 2222 | 2225 | 2232 | rVB  | 457753  | 547209  | 10.21% | 0.642% |
| 95  | 15.374 | 2255 | 2259 | 2265 | rVB  | 1127445 | 1498941 | 27.96% | 1.758% |
| 96  | 15.433 | 2265 | 2269 | 2275 | rBV  | 1635255 | 2133655 | 39.79% | 2.503% |
| 97  | 15.498 | 2276 | 2280 | 2283 | rVV  | 1454942 | 1901690 | 35.47% | 2.230% |
| 98  | 15.527 | 2283 | 2285 | 2292 | rVB2 | 585613  | 799591  | 14.91% | 0.938% |
| 99  | 16.962 | 2524 | 2529 | 2538 | rVB2 | 679351  | 1389878 | 25.92% | 1.630% |
| 100 | 17.403 | 2599 | 2604 | 2615 | rVB  | 502973  | 1023218 | 19.08% | 1.200% |

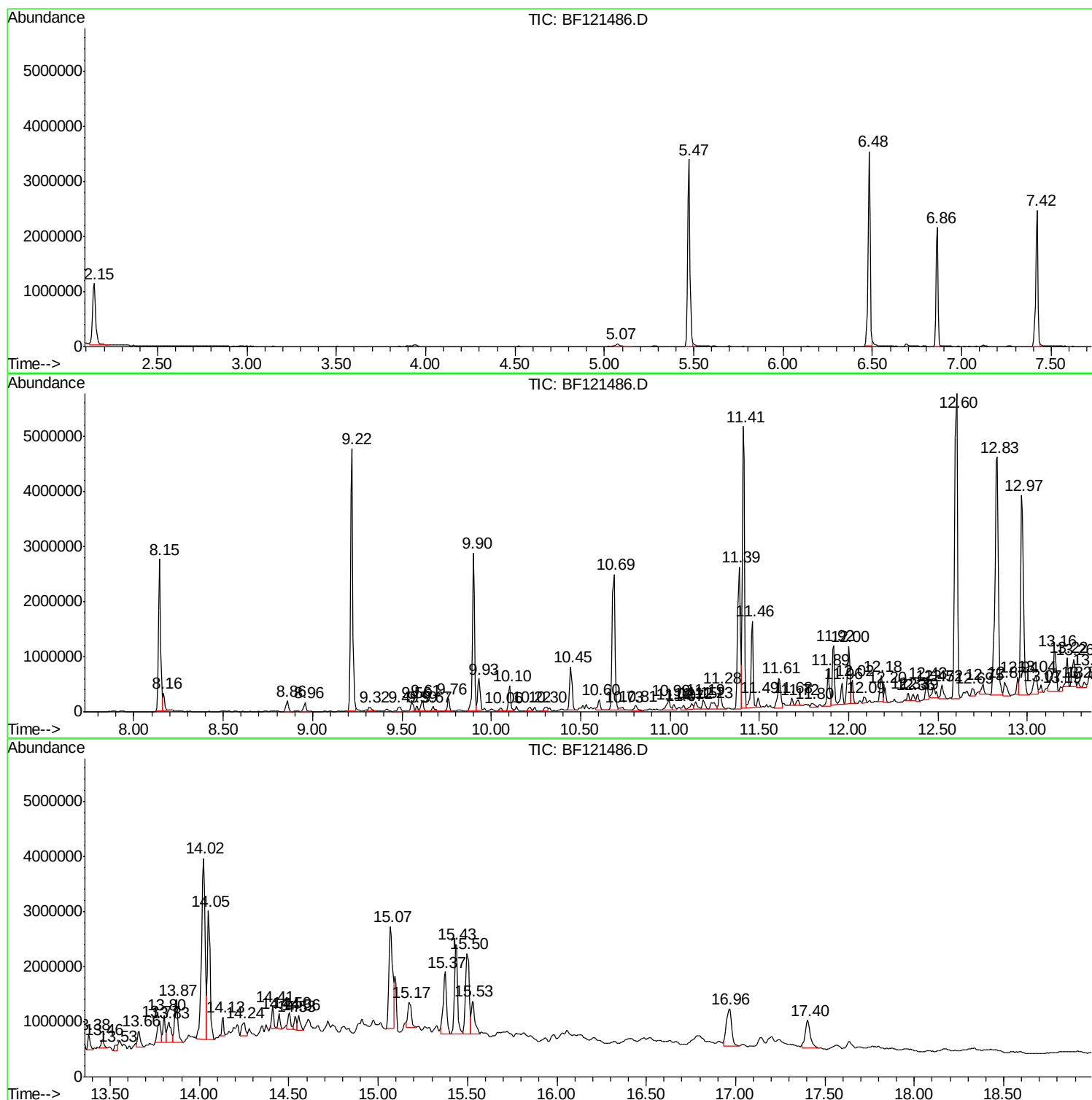
Sum of corrected areas: 85259206

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Instrument :  
BNA\_F  
ClientSampleId :  
LINDEN-JOP-BH-10-B

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
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Instrument :  
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LINDEN-JOP-BH-10-B

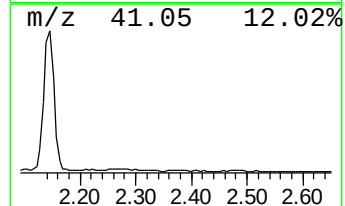
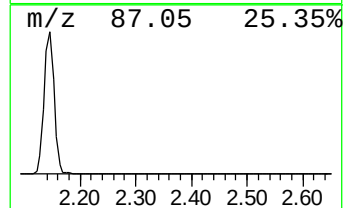
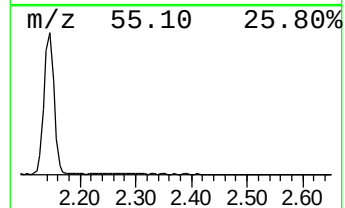
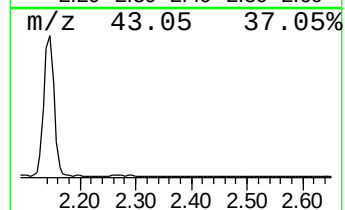
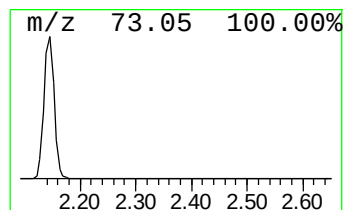
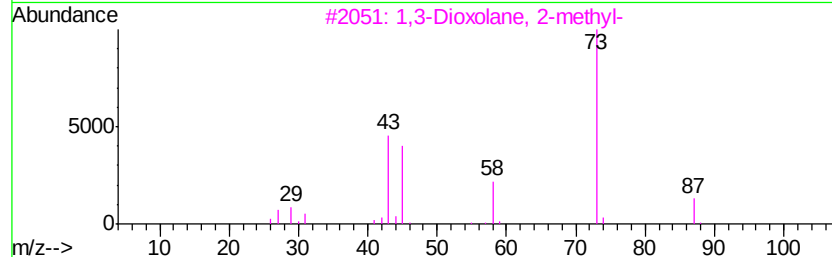
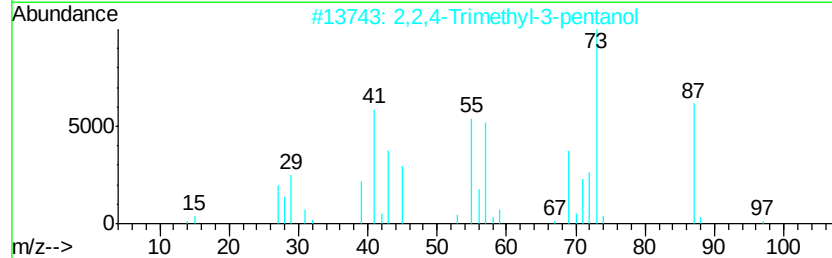
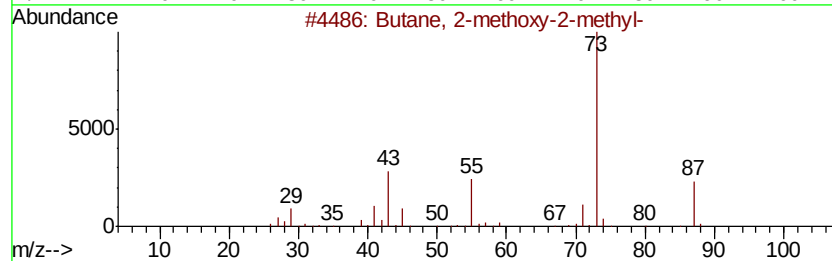
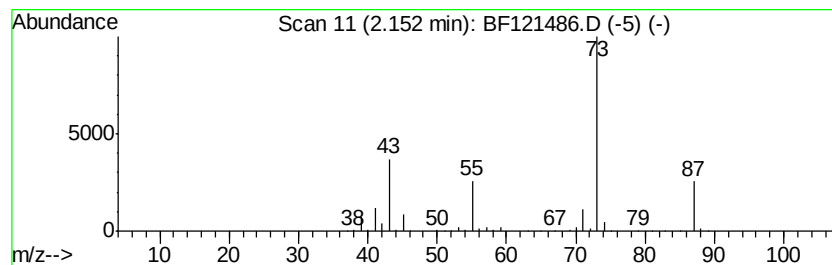
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 2.15 | 15.53 ng | 1431190 | 1,4-Dichlorobenzene-d4 | 6.86 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 37   |
| 4    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 5    |    | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 12   |



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

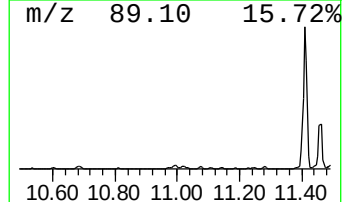
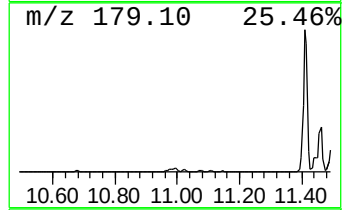
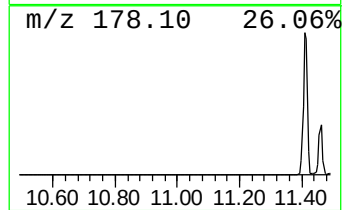
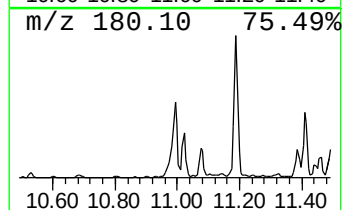
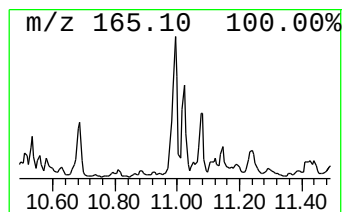
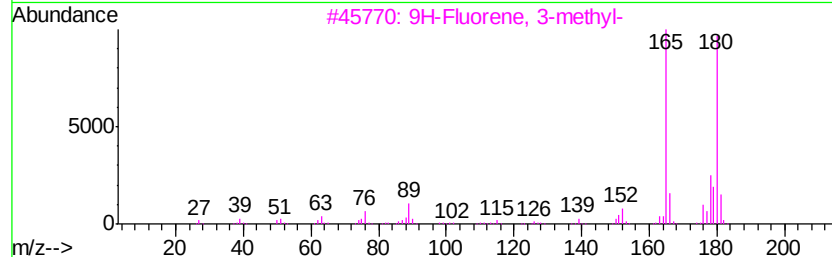
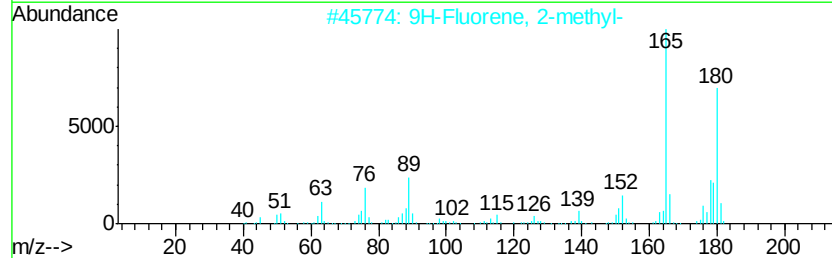
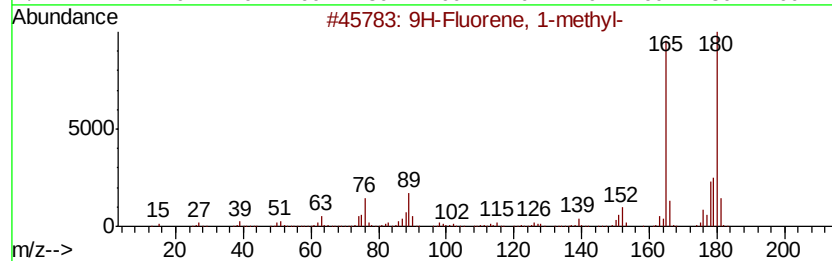
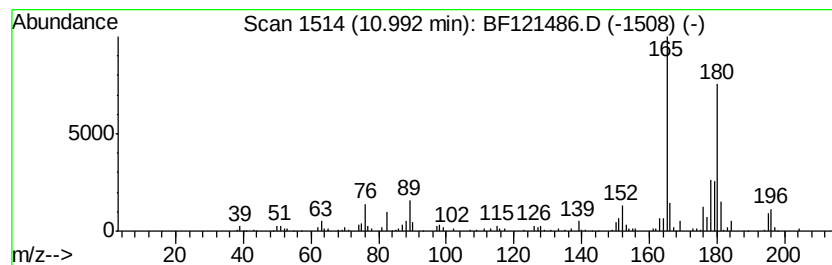
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 9H-Fluorene, 1-methyl- Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.99 | 2.05 ng | 253652 | Phenanthrene-d10 | 11.39 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | 9H-Fluorene, 1-methyl- | 180 | C14H12  | 001730-37-6 | 98   |
| 2    |    | 9H-Fluorene, 2-methyl- | 180 | C14H12  | 001430-97-3 | 97   |
| 3    |    | 9H-Fluorene, 3-methyl- | 180 | C14H12  | 002523-39-9 | 95   |
| 4    |    | 9H-Fluorene, 4-methyl- | 180 | C14H12  | 001556-99-6 | 89   |
| 5    |    | 9H-Fluorene, 9-methyl- | 180 | C14H12  | 002523-37-7 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

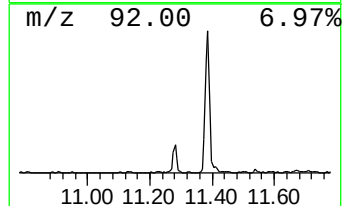
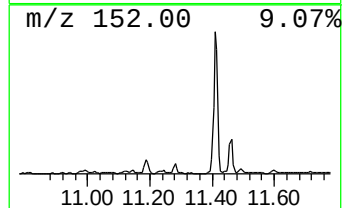
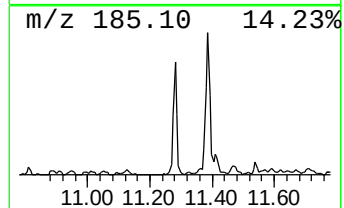
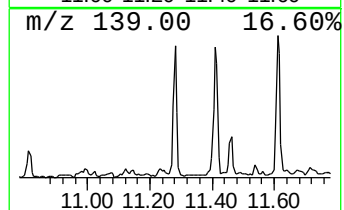
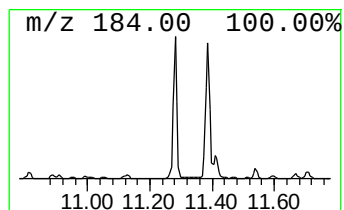
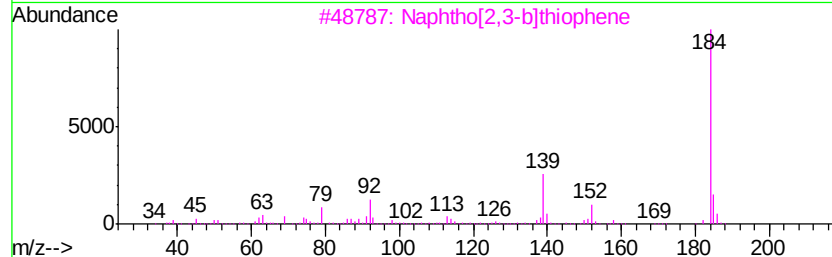
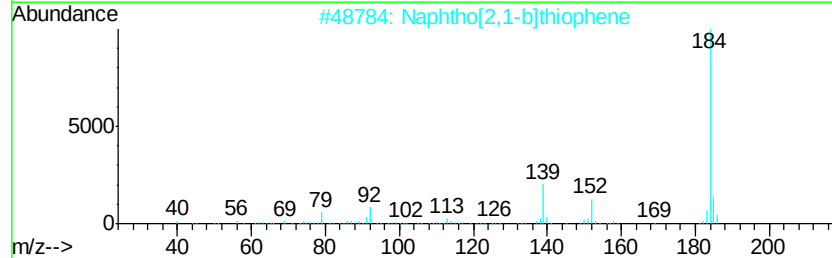
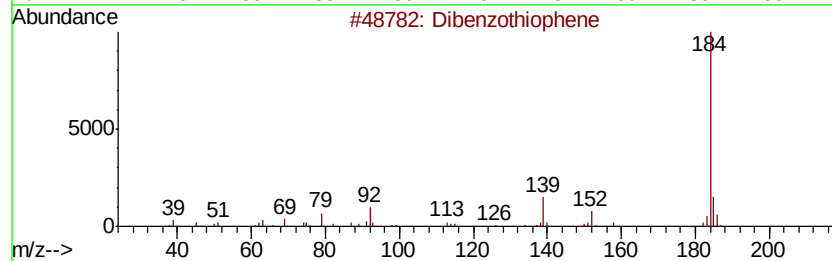
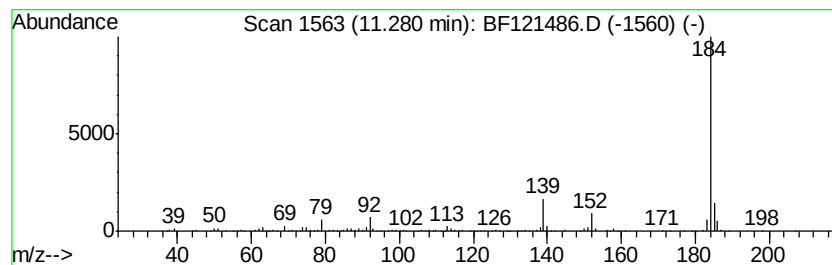
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LINDEN-JOP-BH-10-B

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 3 Dibenzothiophene Concentration Rank 11

| R.T.      | EstConc                 | Area   | Relative to ISTD |             | R.T.  |
|-----------|-------------------------|--------|------------------|-------------|-------|
| 11.28     | 2.81 ng                 | 347729 | Phenanthrene-d10 |             | 11.39 |
| Hit# of 5 | Tentative ID            | MW     | MolForm          | CAS#        | Qual  |
| 1         | Dibenzothiophene        | 184    | C12H8S           | 000132-65-0 | 97    |
| 2         | Naphtho[2,1-b]thiophene | 184    | C12H8S           | 000233-02-3 | 95    |
| 3         | Naphtho[2,3-b]thiophene | 184    | C12H8S           | 000268-77-9 | 95    |
| 4         | Naphtho[1,2-b]thiophene | 184    | C12H8S           | 000234-41-3 | 91    |
| 5         | Azuleno(2,1-b)thiophene | 184    | C12H8S           | 000248-13-5 | 91    |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

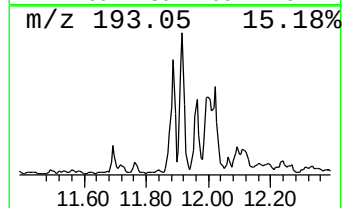
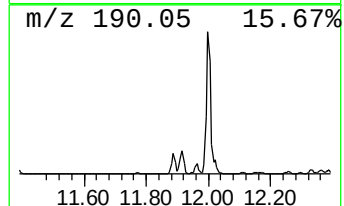
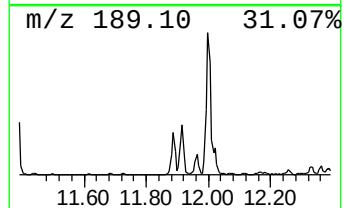
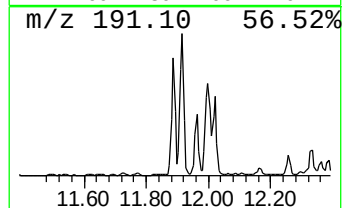
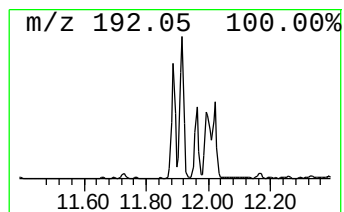
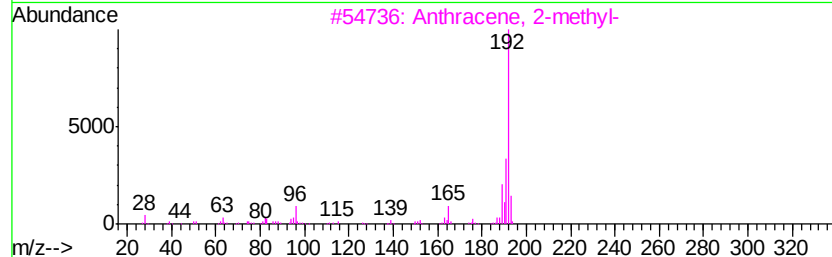
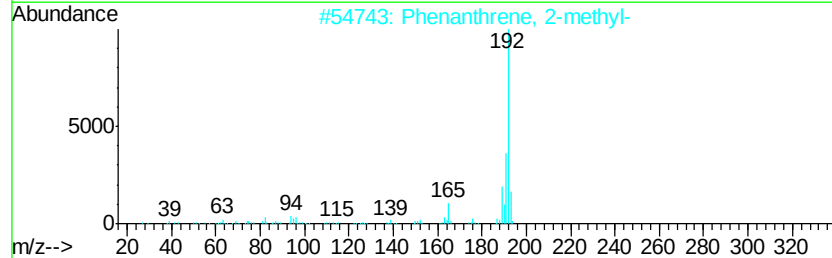
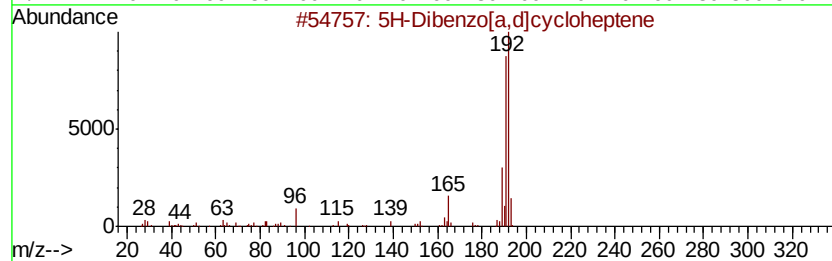
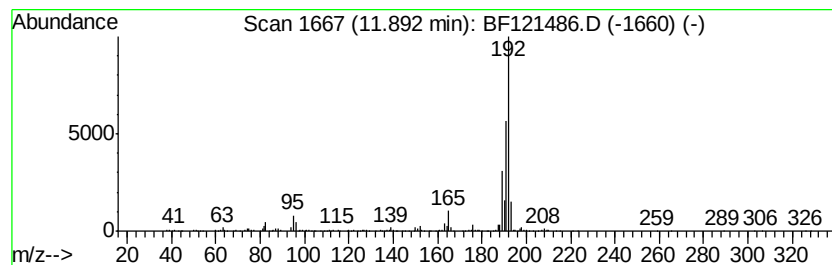
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LINDEN-JOP-BH-10-B

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 4 5H-Dibenzo[a,d]cycloheptene Concentration Rank 6

| R.T.      | EstConc                     | Area   | Relative to ISTD | R.T.        |      |
|-----------|-----------------------------|--------|------------------|-------------|------|
| 11.89     | 4.88 ng                     | 603531 | Phenanthrene-d10 | 11.39       |      |
| Hit# of 5 | Tentative ID                | MW     | MolForm          | CAS#        | Qual |
| 1         | 5H-Dibenzo[a,d]cycloheptene | 192    | C15H12           | 000256-81-5 | 95   |
| 2         | Phenanthrene, 2-methyl-     | 192    | C15H12           | 002531-84-2 | 93   |
| 3         | Anthracene, 2-methyl-       | 192    | C15H12           | 000613-12-7 | 93   |
| 4         | Anthracene, 9-methyl-       | 192    | C15H12           | 000779-02-2 | 91   |
| 5         | Anthracene, 1-methyl-       | 192    | C15H12           | 000610-48-0 | 91   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

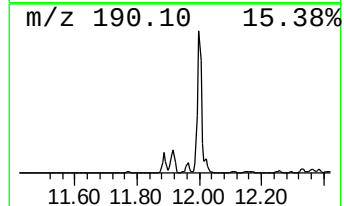
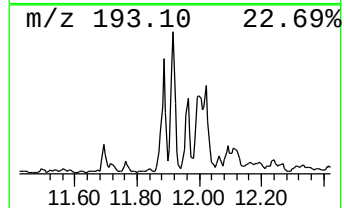
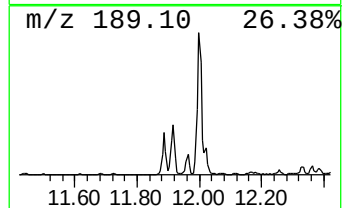
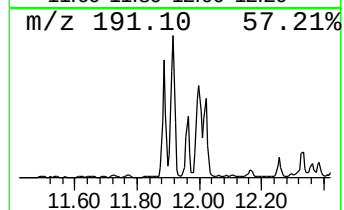
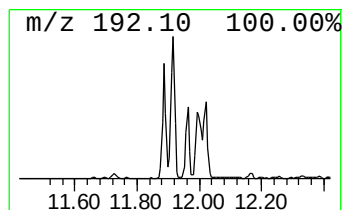
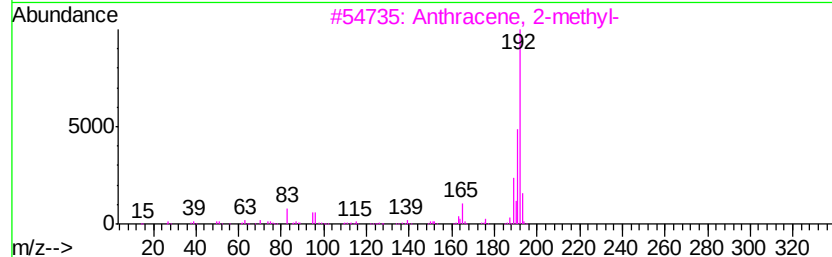
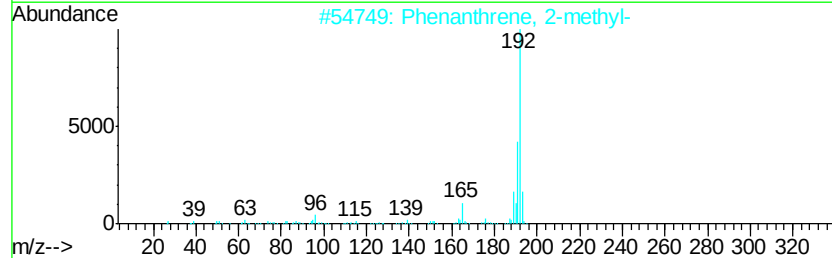
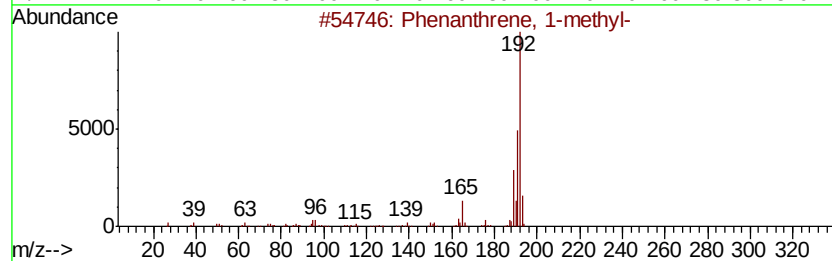
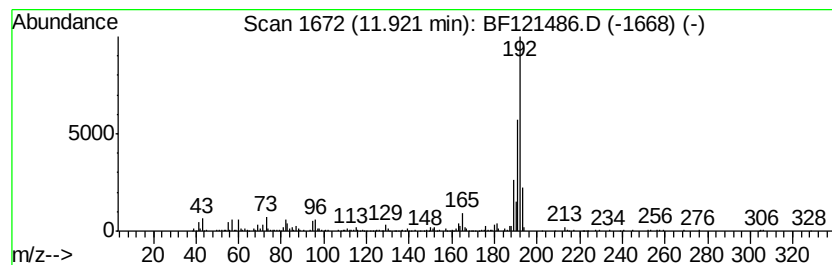
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LINDEN-JOP-BH-10-B

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 5 Phenanthrene, 1-methyl- Concentration Rank 4

| R.T.    | EstConc | Area                    | Relative to ISTD |         | R.T.        |      |
|---------|---------|-------------------------|------------------|---------|-------------|------|
| 11.92   | 8.38 ng | 1036930                 | Phenanthrene-d10 |         | 11.39       |      |
| Hit# of | 5       | Tentative ID            | MW               | MolForm | CAS#        | Qual |
| 1       |         | Phenanthrene, 1-methyl- | 192              | C15H12  | 000832-69-9 | 97   |
| 2       |         | Phenanthrene, 2-methyl- | 192              | C15H12  | 002531-84-2 | 97   |
| 3       |         | Anthracene, 2-methyl-   | 192              | C15H12  | 000613-12-7 | 97   |
| 4       |         | Anthracene, 9-methyl-   | 192              | C15H12  | 000779-02-2 | 94   |
| 5       |         | Anthracene, 1-methyl-   | 192              | C15H12  | 000610-48-0 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
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Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

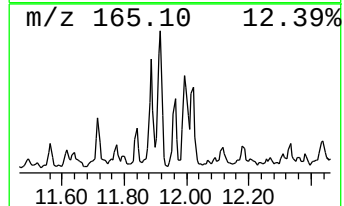
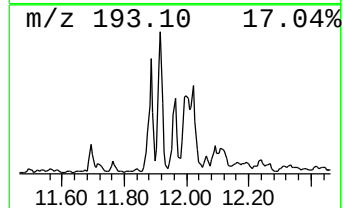
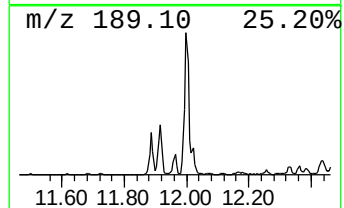
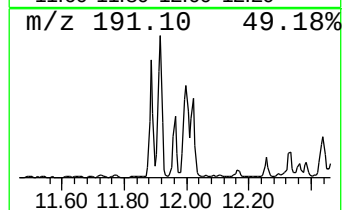
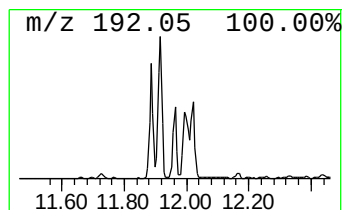
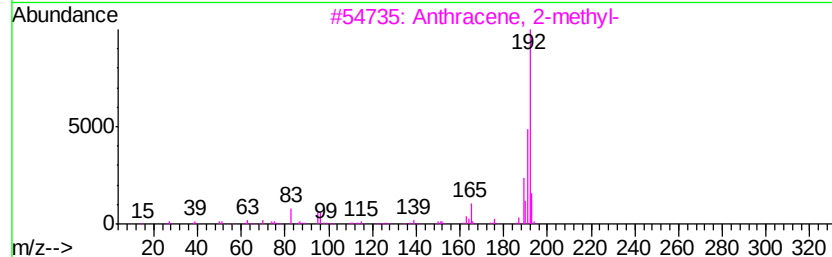
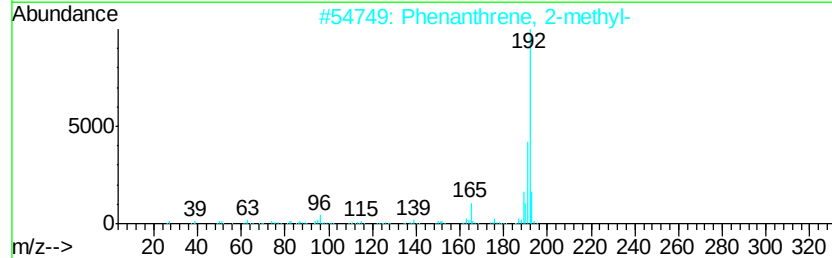
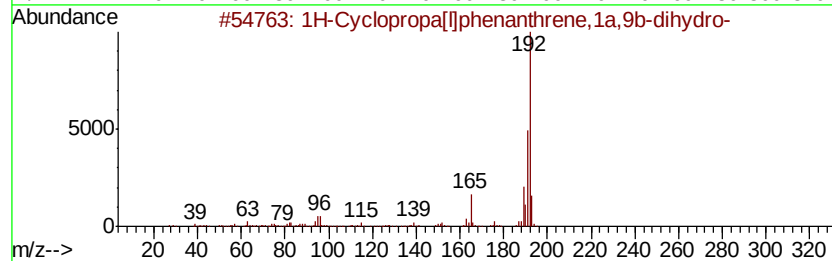
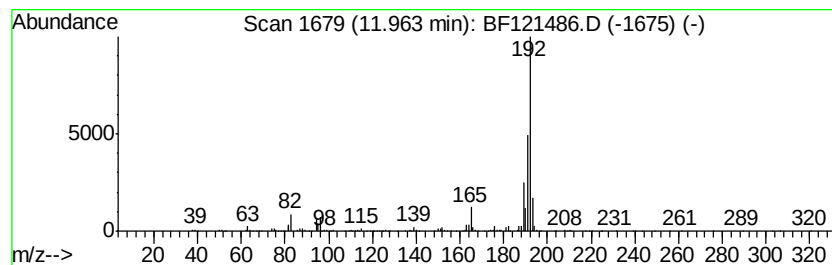
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LINDEN-JOP-BH-10-B

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 6 1H-Cyclopropa[l]phenanthren... Concentration Rank 12

| R.T.      | EstConc                              | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------------|--------|------------------|-------------|------|
| 11.96     | 2.77 ng                              | 342659 | Phenanthrene-d10 | 11.39       |      |
| Hit# of 5 | Tentative ID                         | MW     | MolForm          | CAS#        | Qual |
| 1         | 1H-Cyclopropa[ll]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 96   |
| 2         | Phenanthrene, 2-methyl-              | 192    | C15H12           | 002531-84-2 | 96   |
| 3         | Anthracene, 2-methyl-                | 192    | C15H12           | 000613-12-7 | 96   |
| 4         | 1H-Indene, 1-phenyl-                 | 192    | C15H12           | 001961-96-2 | 94   |
| 5         | Phenanthrene, 1-methyl-              | 192    | C15H12           | 000832-69-9 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LINDEN-JOP-BH-10-B

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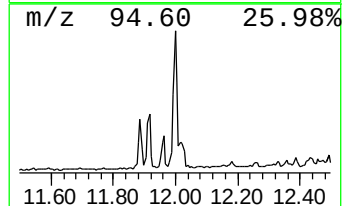
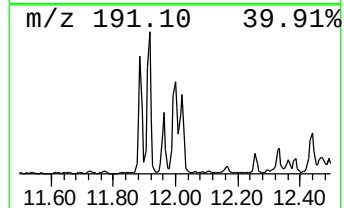
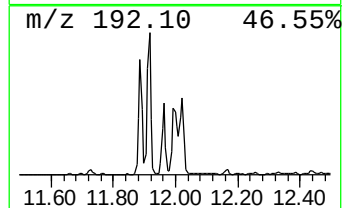
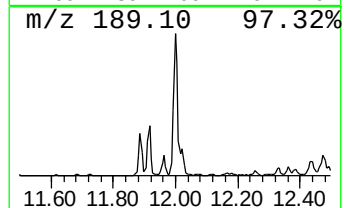
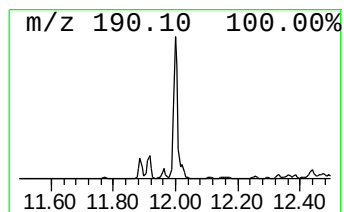
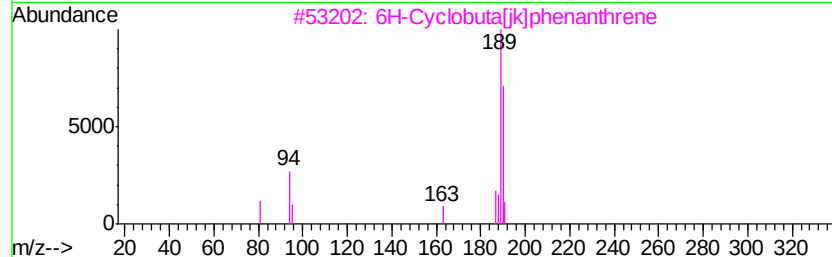
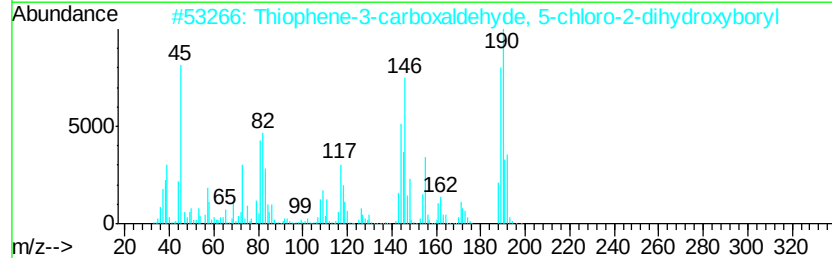
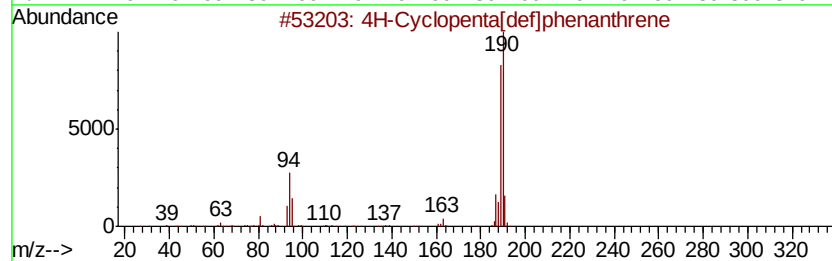
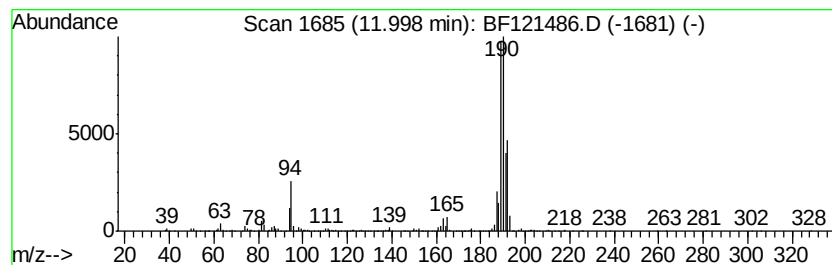
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 12.00 | 8.64 ng | 1068200 | Phenanthrene-d10 | 11.39 |

| Hit# of 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|-----------|-------------------------------------|-----|------------|-------------|------|
| 1         | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10     | 000203-64-5 | 94   |
| 2         | Thiophene-3-carboxaldehyde, 5-ch... | 190 | C5H4BClO3S | 036155-87-0 | 53   |
| 3         | 6H-Cyclobuta[ik]phenanthrene        | 190 | C15H10     | 083469-43-6 | 53   |
| 4         | 2,3,5,6-Tetramethylterephthalald... | 190 | C12H14O2   | 007072-01-7 | 43   |
| 5         | Methyl diselenide                   | 190 | C2H6Se2    | 007101-31-7 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

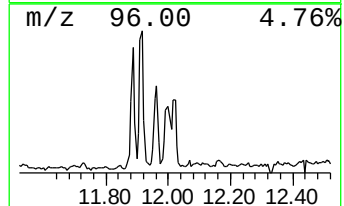
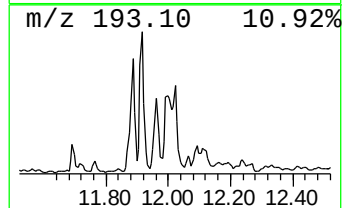
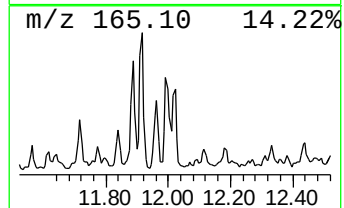
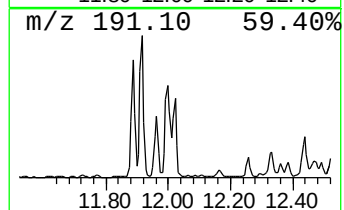
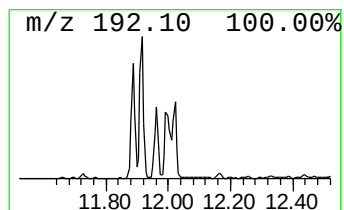
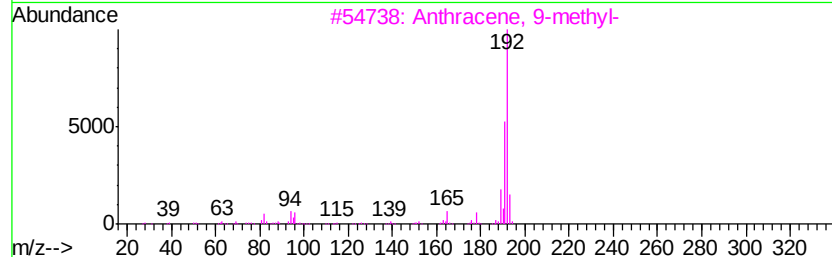
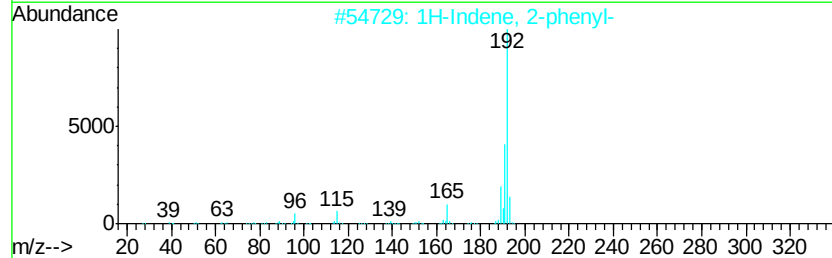
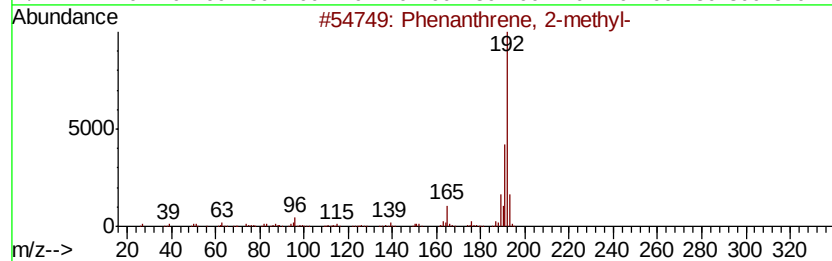
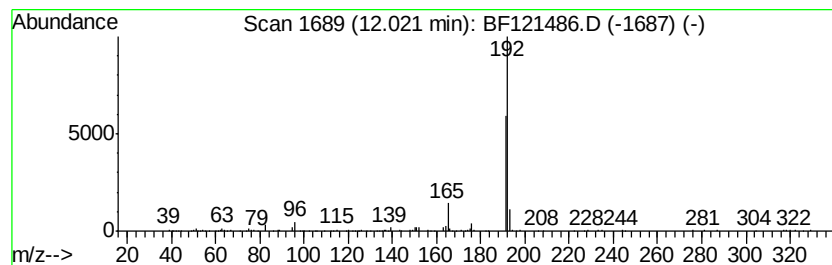
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 8 Phenanthrene, 2-methyl- Concentration Rank 13

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.    |             |      |
|---------|---------|-------------------------|------------------|---------|-------------|------|
| 12.02   | 2.72 ng | 336030                  | Phenanthrene-d10 | 11.39   |             |      |
| Hit# of | 5       | Tentative ID            | MW               | MolForm | CAS#        | Qual |
| 1       |         | Phenanthrene, 2-methyl- | 192              | C15H12  | 002531-84-2 | 87   |
| 2       |         | 1H-Indene, 2-phenyl-    | 192              | C15H12  | 004505-48-0 | 80   |
| 3       |         | Anthracene, 9-methyl-   | 192              | C15H12  | 000779-02-2 | 80   |
| 4       |         | Anthracene, 1-methyl-   | 192              | C15H12  | 000610-48-0 | 80   |
| 5       |         | Phenanthrene, 3-methyl- | 192              | C15H12  | 000832-71-3 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LINDEN-JOP-BH-10-B

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

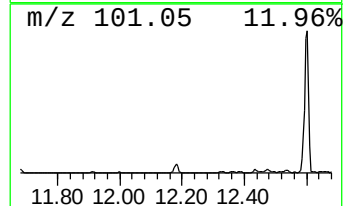
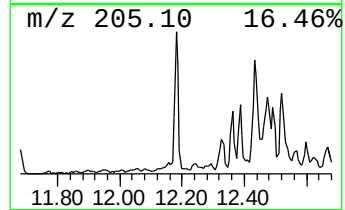
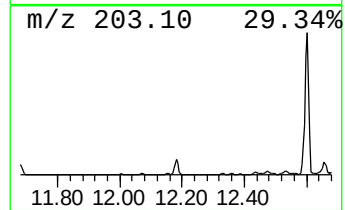
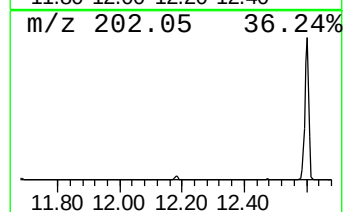
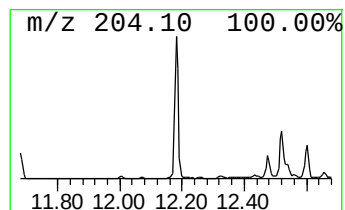
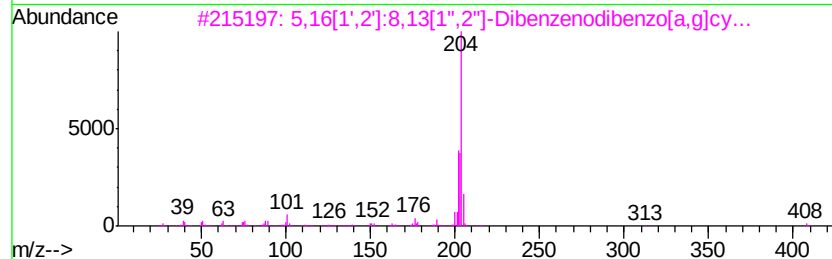
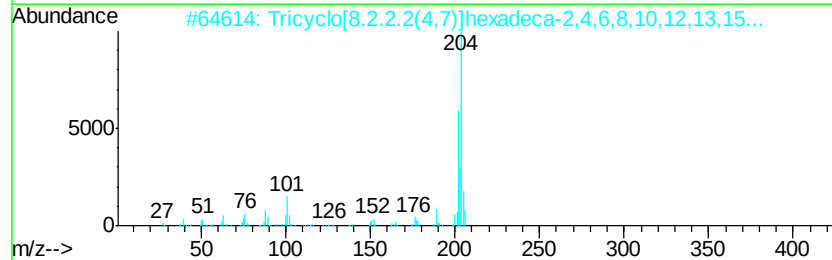
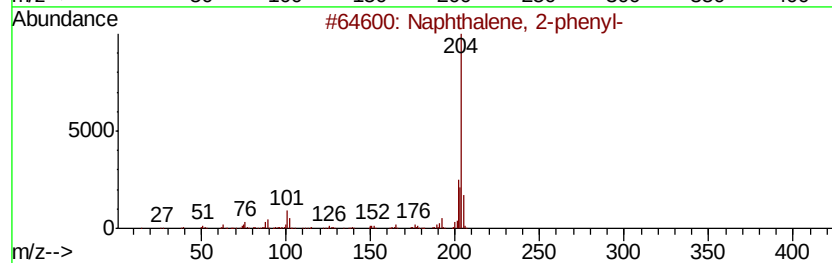
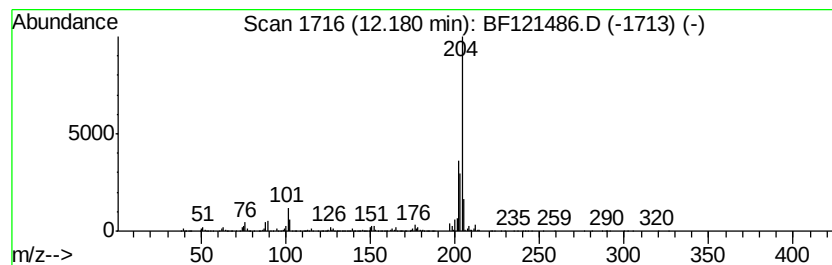
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Naphthalene, 2-phenyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.18 | 3.11 ng | 385143 | Phenanthrene-d10 | 11.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Naphthalene, 2-phenyl-              | 204 | C16H12    | 000612-94-2 | 90   |
| 2    |    | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12    | 006572-60-7 | 72   |
| 3    |    | 5,16[1',2']8,13[1'',2'']-Dibenz...  | 408 | C32H24    | 005672-97-9 | 64   |
| 4    |    | Naphthalene, 1-phenyl-              | 204 | C16H12    | 000605-02-7 | 60   |
| 5    |    | 9,10-Bis(bromomethyl)anthracene     | 362 | C16H12Br2 | 034373-96-1 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

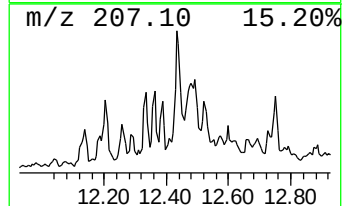
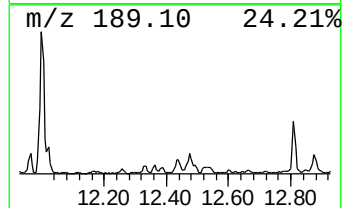
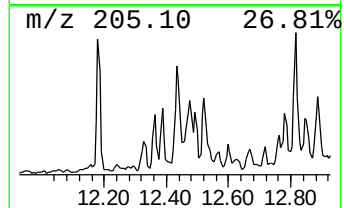
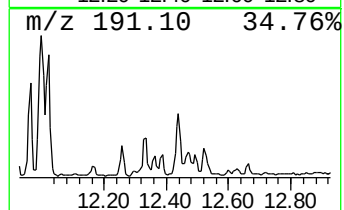
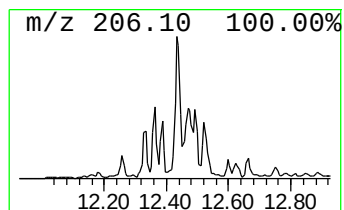
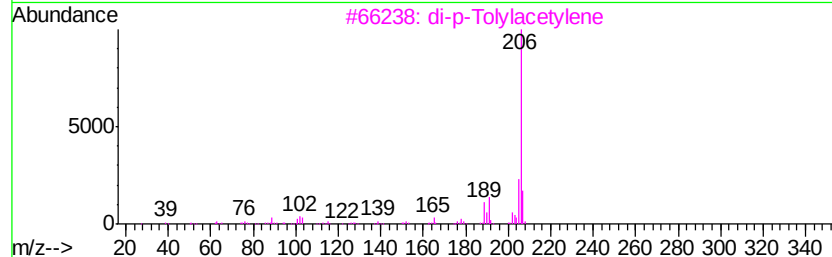
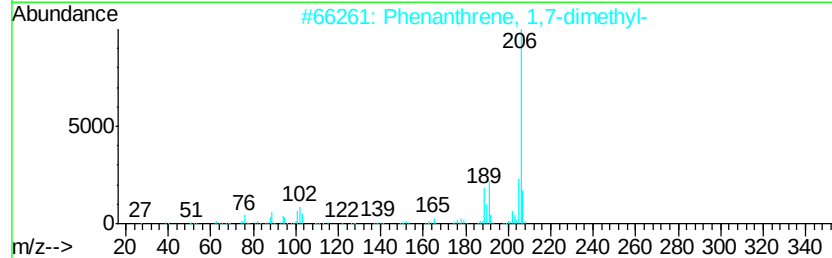
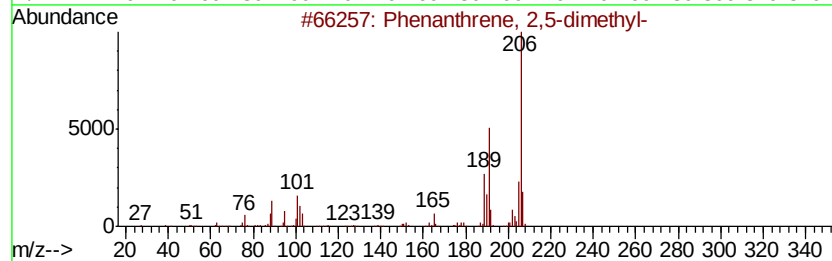
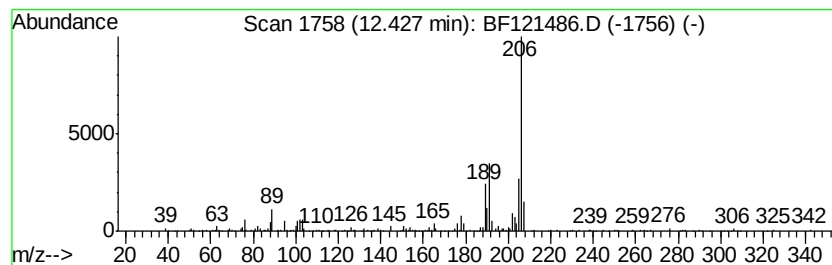
Instrument :  
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 10 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

| R.T.      | EstConc                     | Area   | Relative to ISTD |             | R.T.  |
|-----------|-----------------------------|--------|------------------|-------------|-------|
| 12.43     | 2.83 ng                     | 349563 | Phenanthrene-d10 |             | 11.39 |
| Hit# of 5 | Tentative ID                | MW     | MolForm          | CAS#        | Qual  |
| 1         | Phenanthrene, 2,5-dimethyl- | 206    | C16H14           | 003674-66-6 | 93    |
| 2         | Phenanthrene, 1,7-dimethyl- | 206    | C16H14           | 000483-87-4 | 93    |
| 3         | di-p-Tolylacetylene         | 206    | C16H14           | 002789-88-0 | 93    |
| 4         | Phenanthrene, 3,6-dimethyl- | 206    | C16H14           | 001576-67-6 | 93    |
| 5         | 9,10-Dimethylantracene      | 206    | C16H14           | 000781-43-1 | 91    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LINDEN-JOP-BH-10-B

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

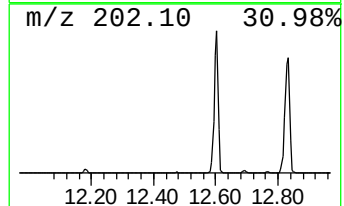
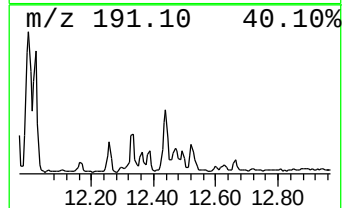
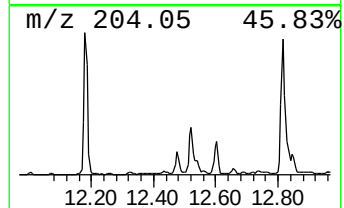
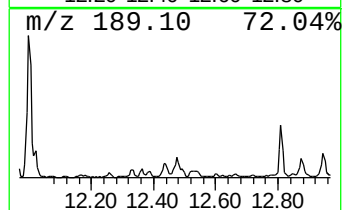
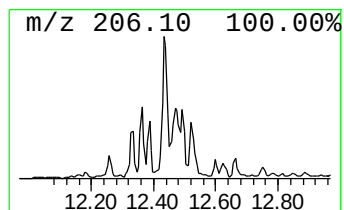
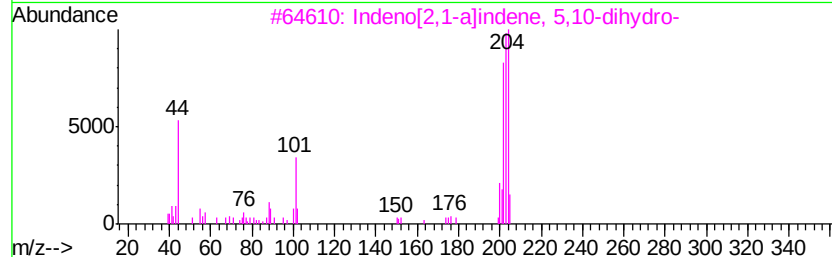
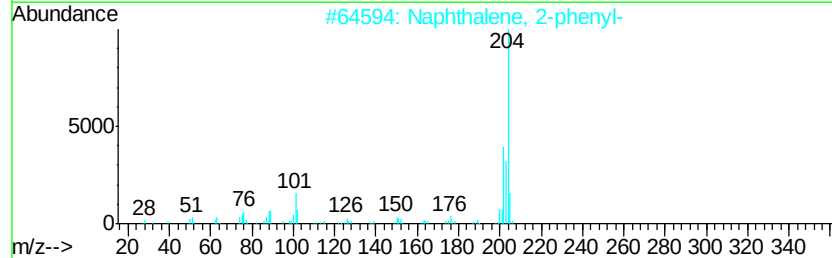
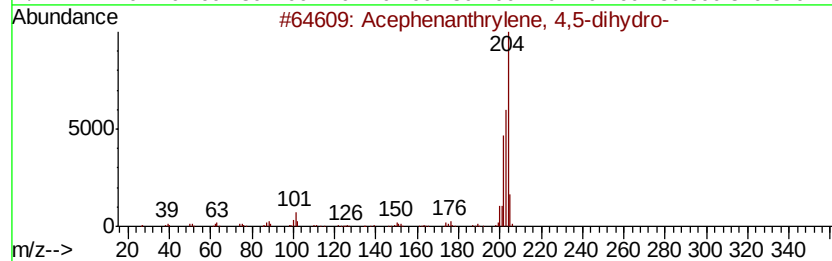
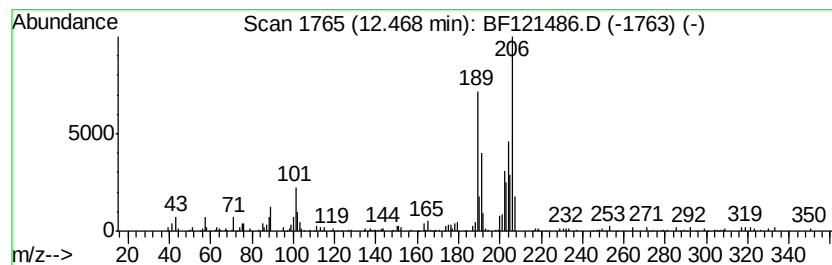
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 unknown12.47 Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.47 | 2.47 ng | 305312 | Phenanthrene-d10 | 11.39 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Acephenanthrylene, 4,5-dihydro-      | 204 | C16H12  | 006232-48-0 | 49   |
| 2    |    | Naphthalene, 2-phenyl-               | 204 | C16H12  | 000612-94-2 | 42   |
| 3    |    | Indeno[2,1-a]indene, 5,10-dihydro-   | 204 | C16H12  | 006543-29-9 | 35   |
| 4    |    | Naphthalene, 1-phenyl-               | 204 | C16H12  | 000605-02-7 | 35   |
| 5    |    | Benzene, 1,1'-(1,3-butadienyldide... | 206 | C16H14  | 004165-81-5 | 25   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LINDEN-JOP-BH-10-B

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

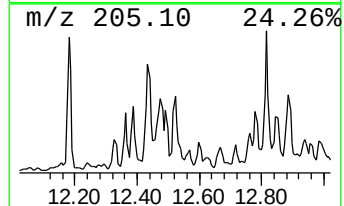
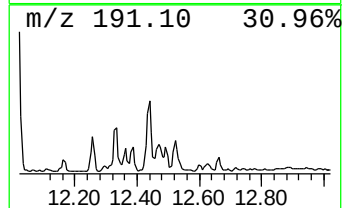
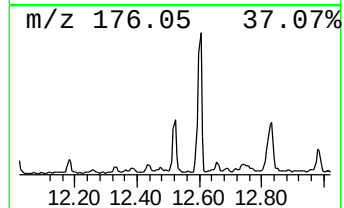
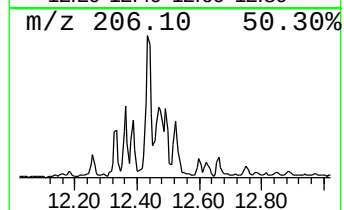
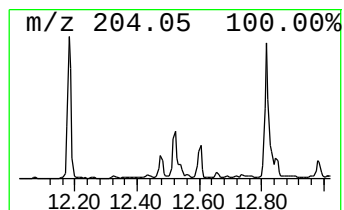
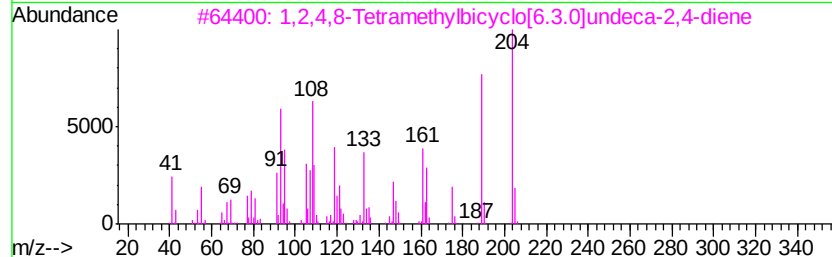
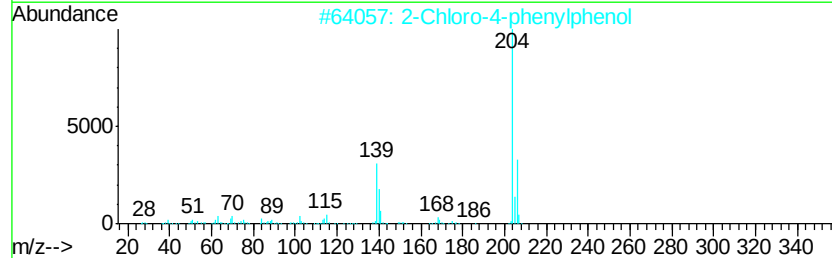
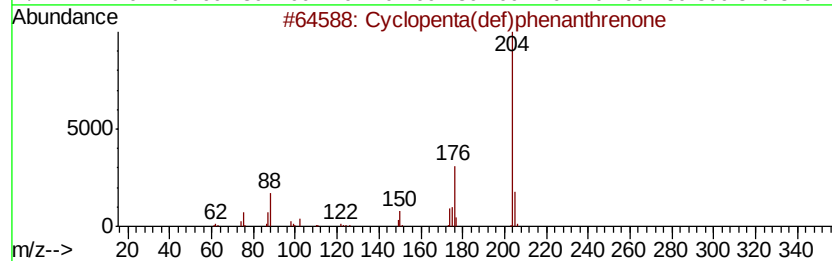
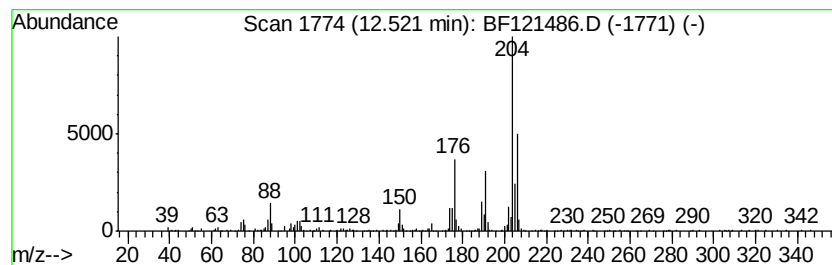
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Cyclopenta(def)phenanthrenone Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.52 | 2.33 ng | 288097 | Phenanthrene-d10 | 11.39 |

| Hit# | of | Tentative ID                                      | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrenone                     | 204 | C15H8O   | 005737-13-3 | 91   |
| 2    |    | 2-Chloro-4-phenylphenol                           | 204 | C12H9ClO | 000092-04-6 | 46   |
| 3    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene | 204 | C15H24   | 137235-51-9 | 41   |
| 4    |    | Phenanthrene, 3,6-dimethyl-                       | 206 | C16H14   | 001576-67-6 | 38   |
| 5    |    | [14]Annulene, 1,6:8,13-bis(metha...               | 206 | C16H14   | 085385-68-8 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

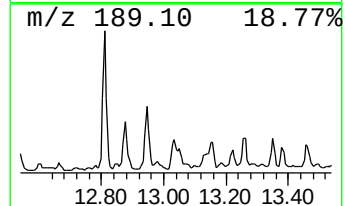
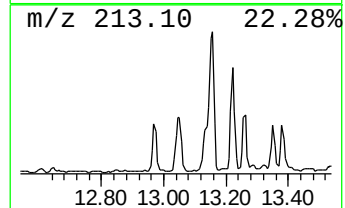
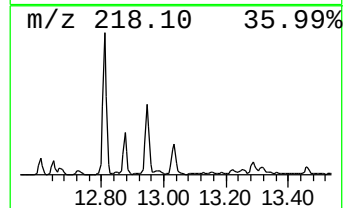
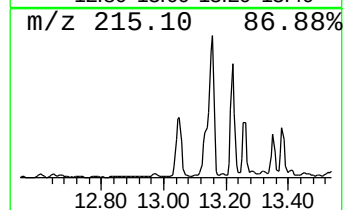
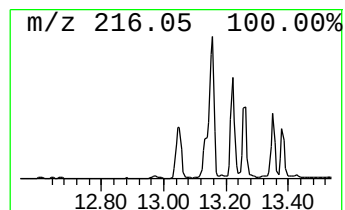
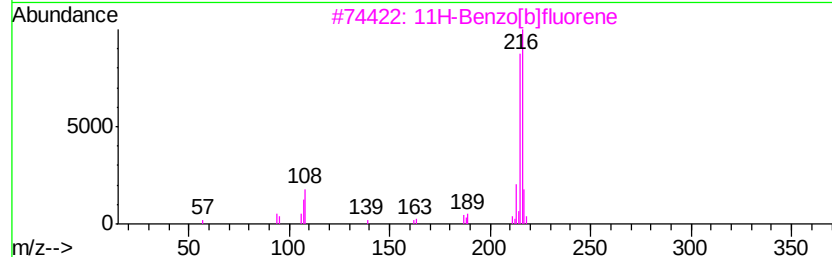
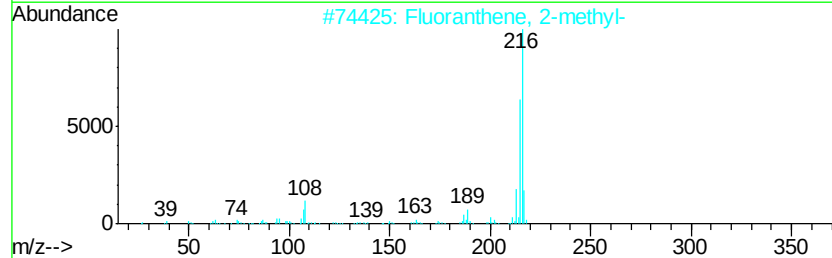
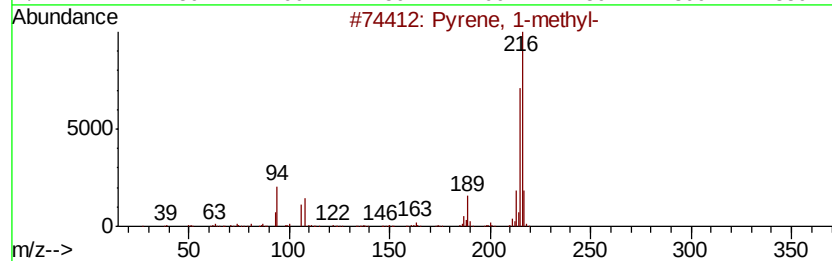
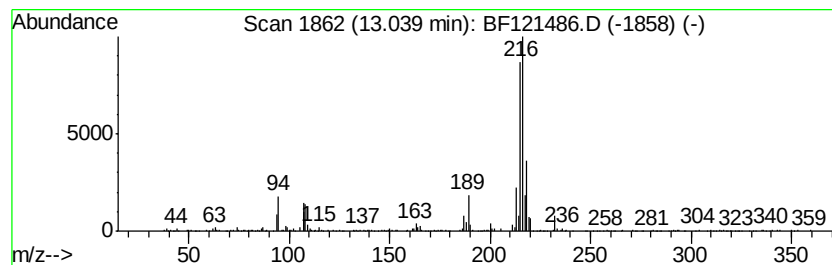
Instrument :  
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 13 Pyrene, 1-methyl- Concentration Rank 16

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.           |
|---------|---------|-------------------------|------------------|----------------|
| 13.04   | 2.31 ng | 558462                  | Chrysene-d12     | 14.03          |
| Hit# of | 5       | Tentative ID            | MW MolForm       | CAS# Qual      |
| 1       |         | Pyrene, 1-methyl-       | 216 C17H12       | 002381-21-7 96 |
| 2       |         | Fluoranthene, 2-methyl- | 216 C17H12       | 033543-31-6 96 |
| 3       |         | 11H-Benzo[b]fluorene    | 216 C17H12       | 000243-17-4 94 |
| 4       |         | Pyrene, 4-methyl-       | 216 C17H12       | 003353-12-6 90 |
| 5       |         | 11H-Benzo[a]fluorene    | 216 C17H12       | 000238-84-6 74 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

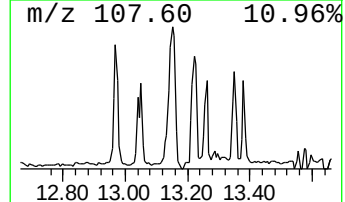
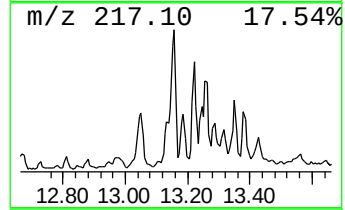
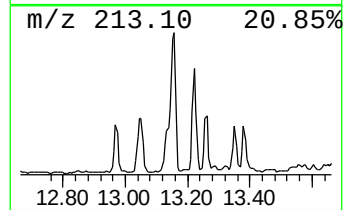
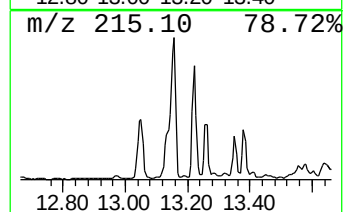
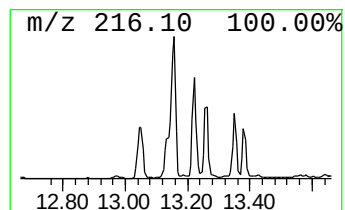
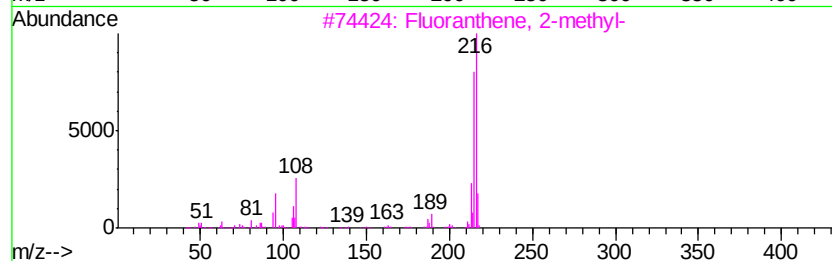
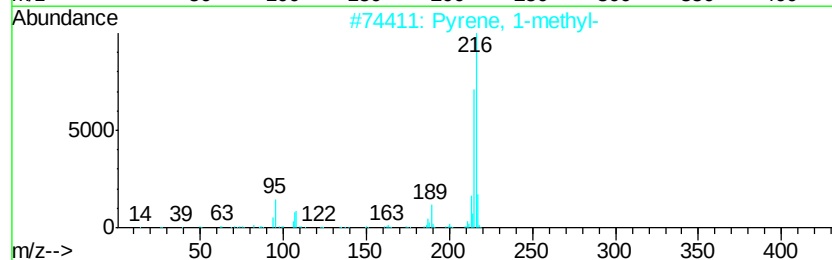
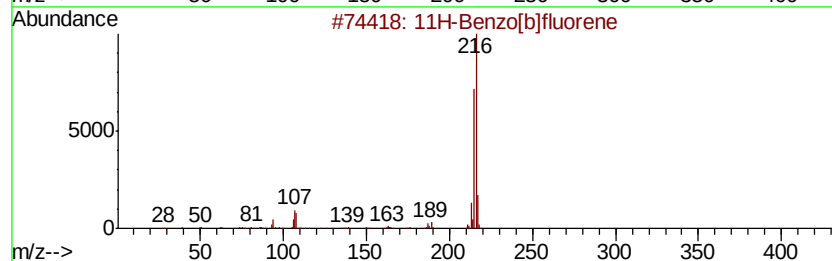
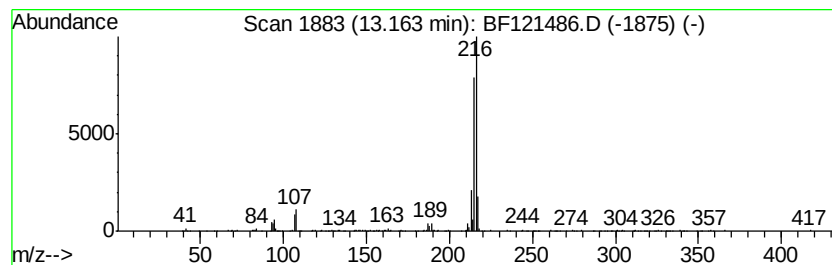
Instrument :  
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 11H-Benzo[b]fluorene Concentration Rank 7

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.           |
|---------|---------|-------------------------|------------------|----------------|
| 13.16   | 4.16 ng | 1002850                 | Chrysene-d12     | 14.03          |
| Hit# of | 5       | Tentative ID            | MW MolForm       | CAS# Qual      |
| 1       |         | 11H-Benzo[b]fluorene    | 216 C17H12       | 000243-17-4 93 |
| 2       |         | Pyrene, 1-methyl-       | 216 C17H12       | 002381-21-7 93 |
| 3       |         | Fluoranthene, 2-methyl- | 216 C17H12       | 033543-31-6 91 |
| 4       |         | 11H-Benzo[a]fluorene    | 216 C17H12       | 000238-84-6 91 |
| 5       |         | 7H-Benzo[c]fluorene     | 216 C17H12       | 000205-12-9 72 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

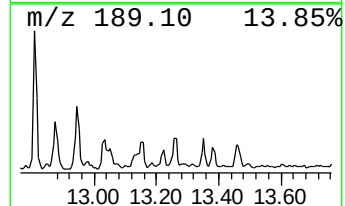
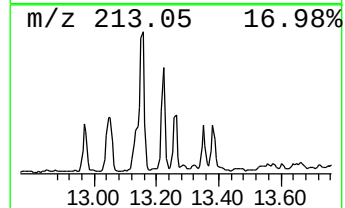
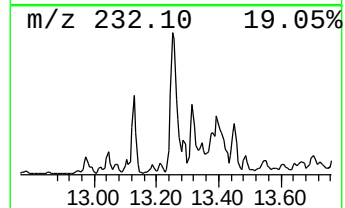
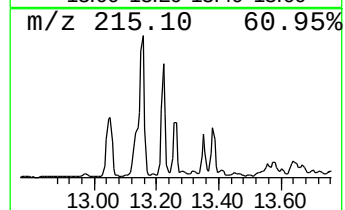
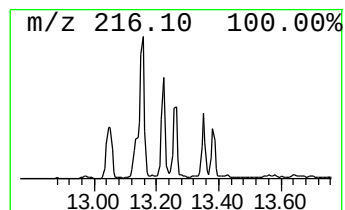
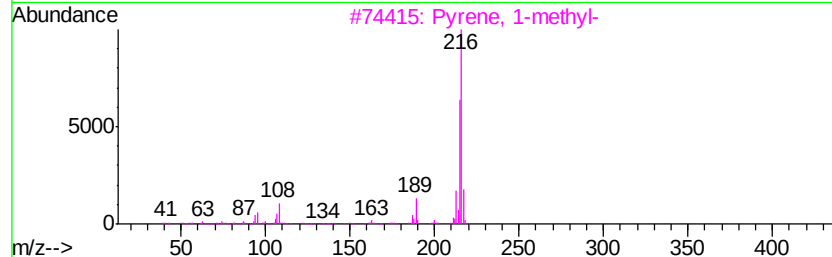
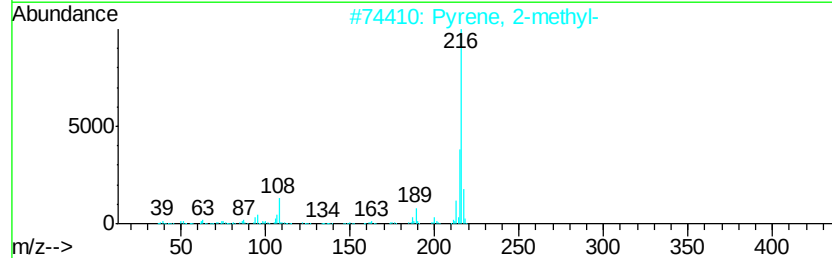
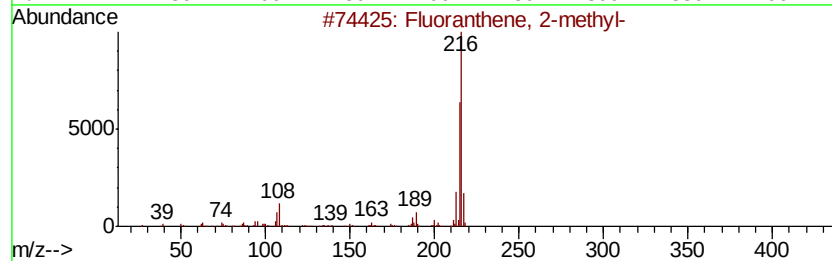
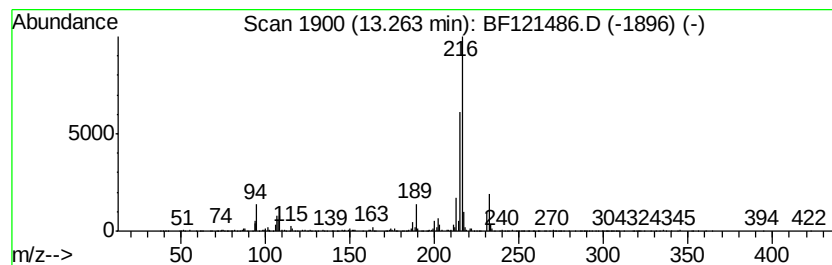
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 Fluoranthene, 2-methyl- Concentration Rank 20

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.           |
|---------|---------|-------------------------|------------------|----------------|
| 13.26   | 2.03 ng | 490474                  | Chrysene-d12     | 14.03          |
| Hit# of | 5       | Tentative ID            | MW MolForm       | CAS# Qual      |
| 1       |         | Fluoranthene, 2-methyl- | 216 C17H12       | 033543-31-6 78 |
| 2       |         | Pyrene, 2-methyl-       | 216 C17H12       | 003442-78-2 78 |
| 3       |         | Pyrene, 1-methyl-       | 216 C17H12       | 002381-21-7 70 |
| 4       |         | 11H-Benzo[alfluorene    | 216 C17H12       | 000238-84-6 60 |
| 5       |         | Pyrene, 4-methyl-       | 216 C17H12       | 003353-12-6 49 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

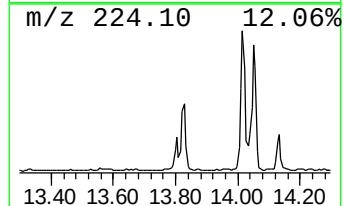
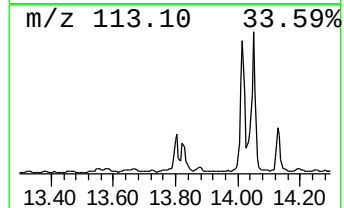
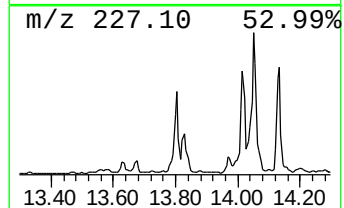
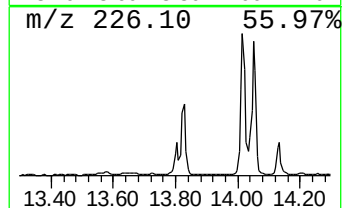
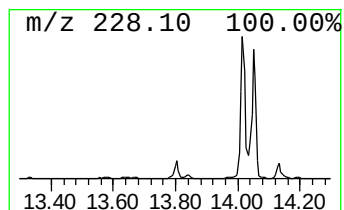
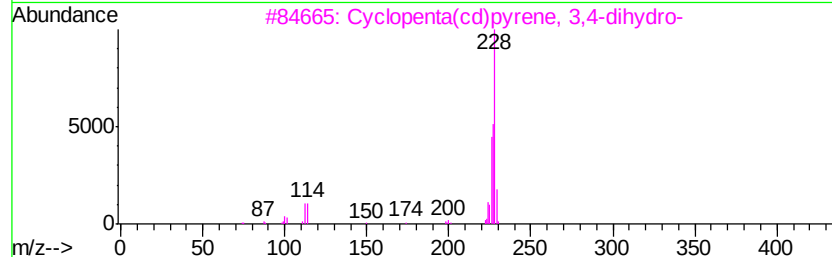
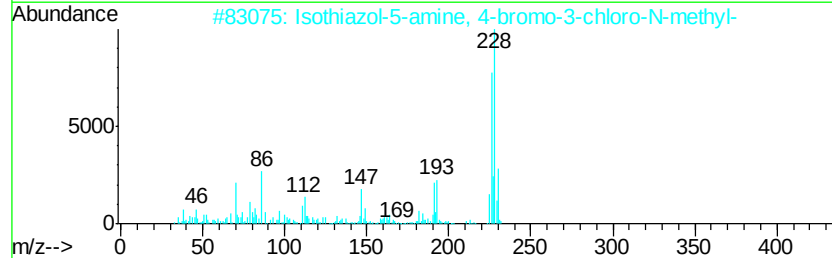
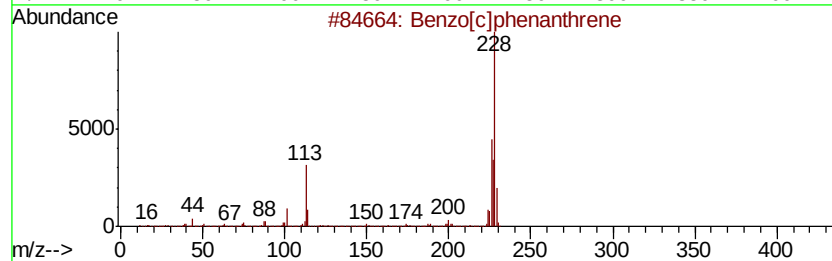
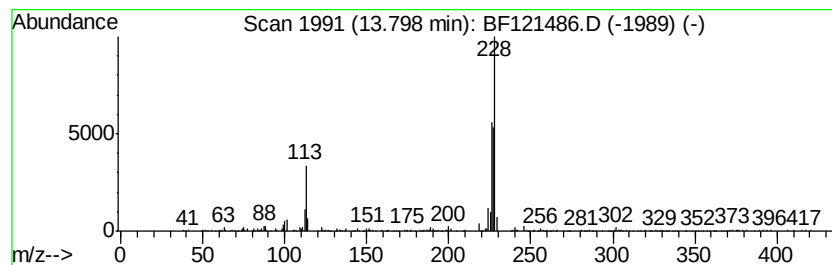
Instrument :  
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 16 Benzo[c]phenanthrene Concentration Rank 18

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.        |              |      |
|---------|---------|-------------------------------------|------------------|-------------|--------------|------|
| 13.80   | 2.06 ng | 498115                              | Chrysene-d12     | 14.03       |              |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm     | CAS#         | Qual |
| 1       |         | Benzo[c]phenanthrene                | 228              | C18H12      | 000195-19-7  | 53   |
| 2       |         | Isothiazol-5-amine, 4-bromo-3-ch... | 226              | C4H4BrClN2S | 1000272-59-9 | 50   |
| 3       |         | Cyclopenta(cd)pyrene, 3,4-dihydro-  | 228              | C18H12      | 025732-74-5  | 49   |
| 4       |         | Triphenylene                        | 228              | C18H12      | 000217-59-4  | 43   |
| 5       |         | Benz[a]anthracene                   | 228              | C18H12      | 000056-55-3  | 41   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

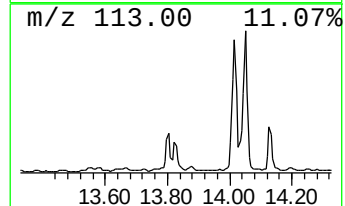
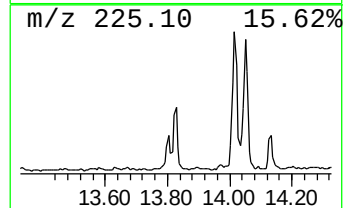
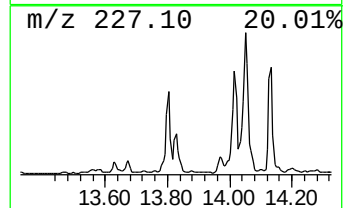
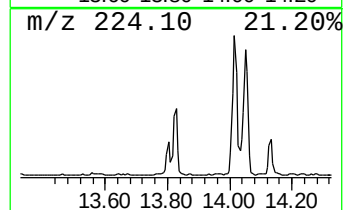
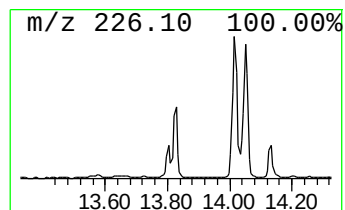
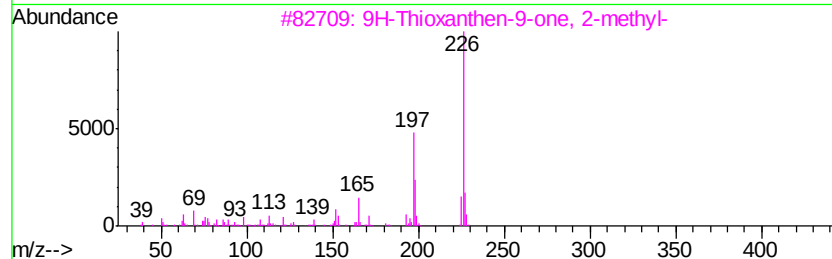
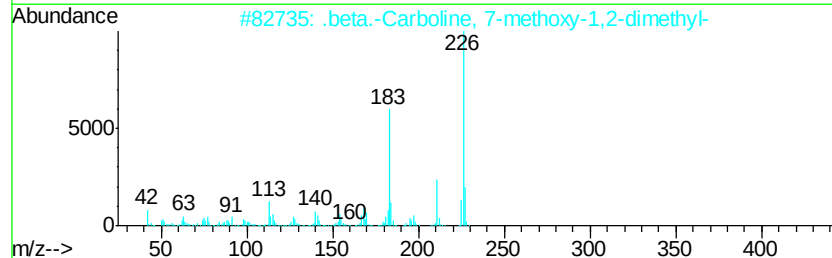
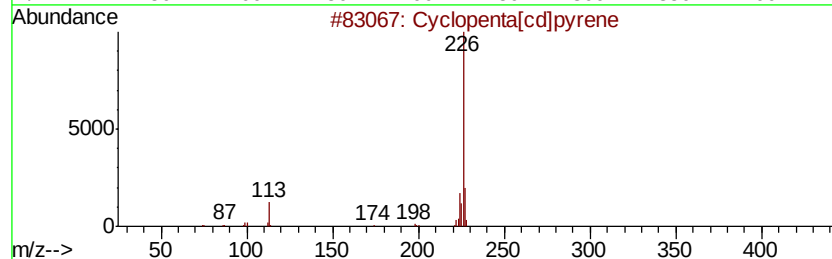
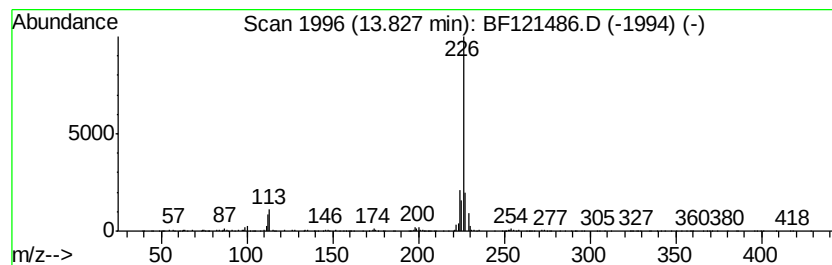
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 Cyclopenta[cd]pyrene Concentration Rank 17

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.           |
|---------|---------|-------------------------------------|------------------|----------------|
| 13.83   | 2.20 ng | 531560                              | Chrysene-d12     | 14.03          |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |         | Cyclopenta[cd]pyrene                | 226 C18H10       | 027208-37-3 93 |
| 2       |         | .beta.-Carboline, 7-methoxy-1,2-... | 226 C14H14N2O    | 006519-18-2 58 |
| 3       |         | 9H-Thioxanthen-9-one, 2-methyl-     | 226 C14H10OS     | 015774-82-0 53 |
| 4       |         | Benzene, 1-bromo-2,6-dichloro-      | 224 C6H3BrCl2    | 019393-92-1 42 |
| 5       |         | 1,8-Diazacyclotetradecane-2,7-dione | 226 C12H22N2O2   | 004266-66-4 38 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

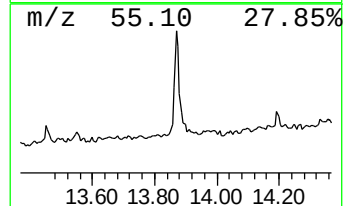
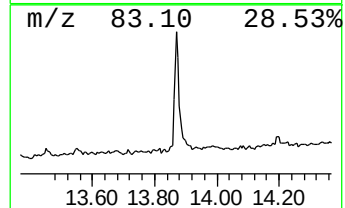
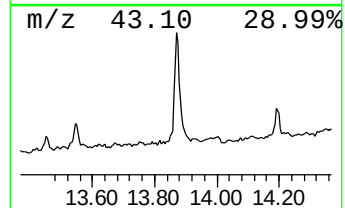
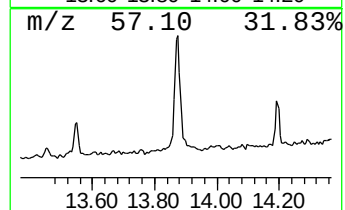
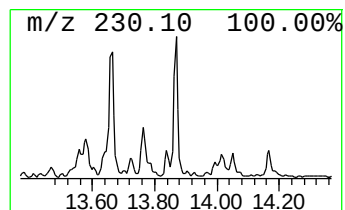
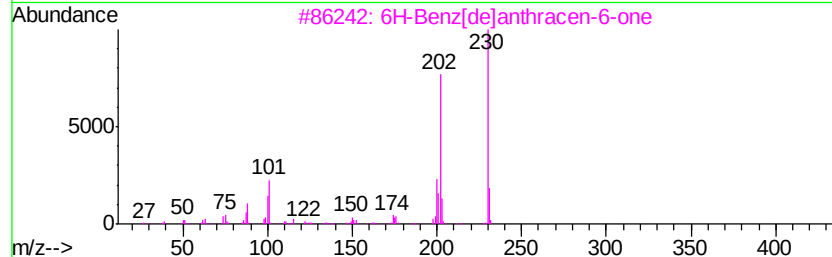
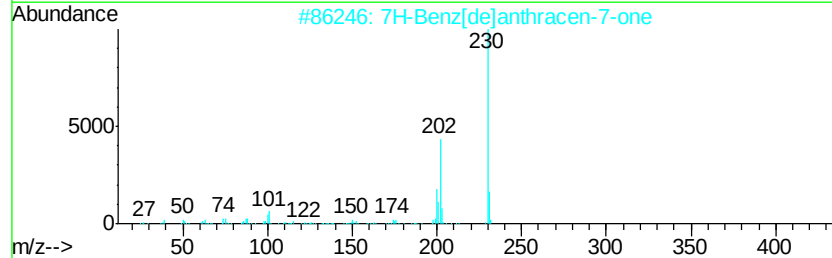
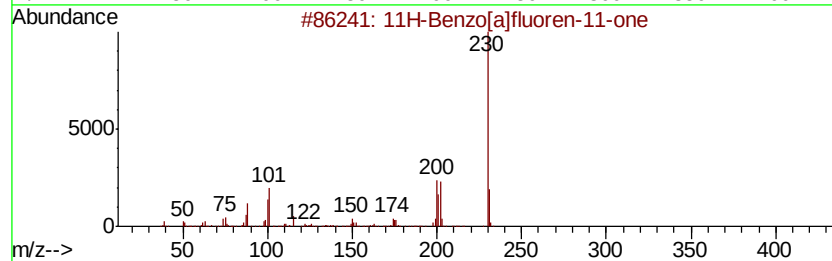
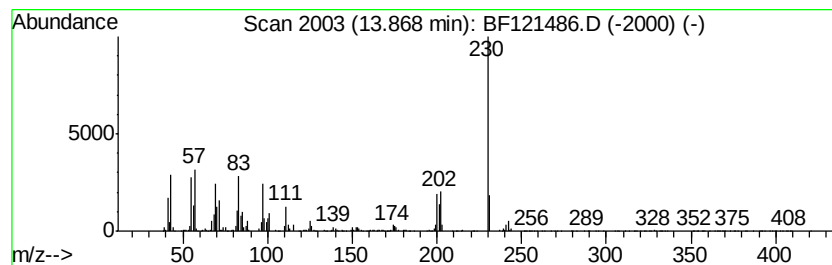
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 18 11H-Benzo[a]fluoren-11-one Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD           |     | R.T.    |             |      |
|-------|---------|--------|----------------------------|-----|---------|-------------|------|
| 13.87 | 3.64 ng | 879208 | Chrysene-d12               |     | 14.03   |             |      |
| Hit#  | of      | 5      | Tentative ID               | MW  | MolForm | CAS#        | Qual |
| 1     |         |        | 11H-Benzo[a]fluoren-11-one | 230 | C17H10O | 000479-79-8 | 95   |
| 2     |         |        | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 78   |
| 3     |         |        | 6H-Benz[de]anthracen-6-one | 230 | C17H10O | 080252-14-8 | 53   |
| 4     |         |        | m-Terphenyl                | 230 | C18H14  | 000092-06-8 | 38   |
| 5     |         |        | o-Terphenyl                | 230 | C18H14  | 000084-15-1 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

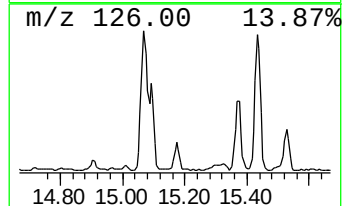
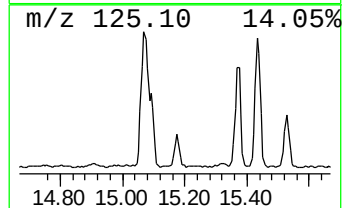
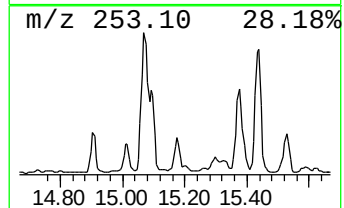
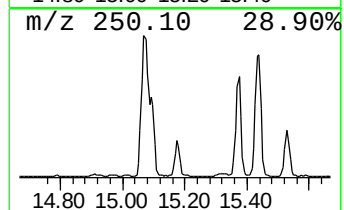
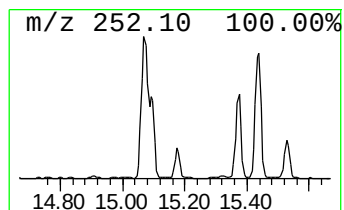
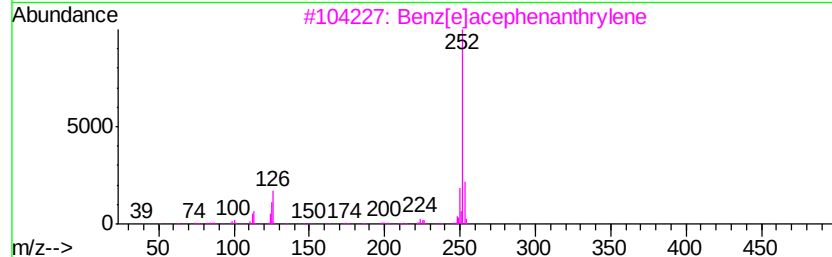
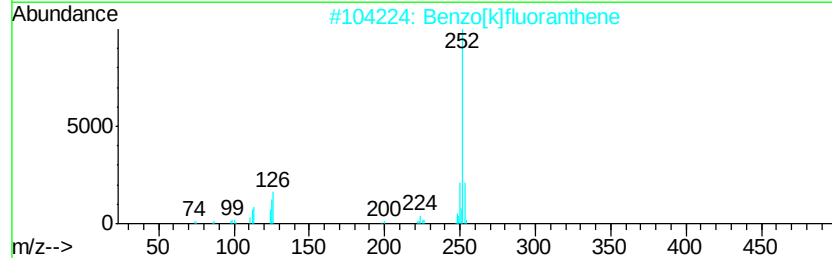
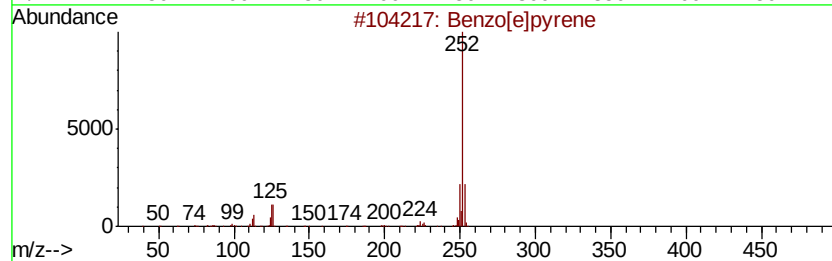
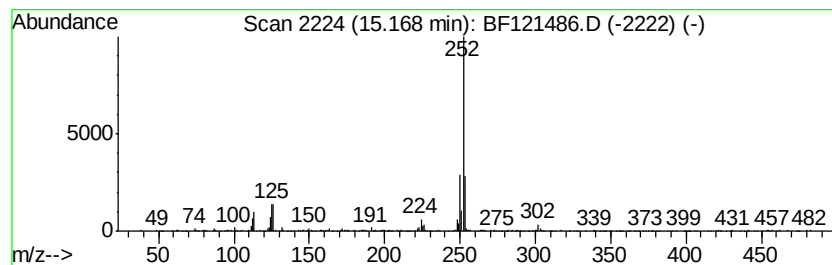
Instrument :  
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ClientSampleId :  
LINDEN-JOP-BH-10-B

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 19 Benzo[e]pyrene Concentration Rank 5

| R.T.      | EstConc                   | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------------|--------|------------------|-------------|------|
| 15.17     | 5.75 ng                   | 547209 | Perylene-d12     | 15.50       |      |
| Hit# of 5 | Tentative ID              | MW     | MolForm          | CAS#        | Qual |
| 1         | Benzo[e]pyrene            | 252    | C20H12           | 000192-97-2 | 96   |
| 2         | Benzo[k]fluoranthene      | 252    | C20H12           | 000207-08-9 | 96   |
| 3         | Benzo[e]acephenanthrylene | 252    | C20H12           | 000205-99-2 | 93   |
| 4         | Benzo[a]pyrene            | 252    | C20H12           | 000050-32-8 | 91   |
| 5         | Perylene                  | 252    | C20H12           | 000198-55-0 | 81   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083120\  
Data File : BF121486.D  
Acq On : 31 Aug 2020 19:32  
Operator : JU/CG  
Sample : L3847-15  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
LINDEN-JOP-BH-10-B

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

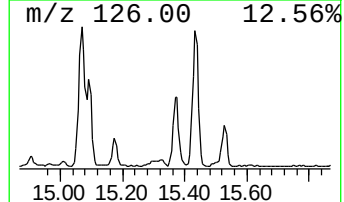
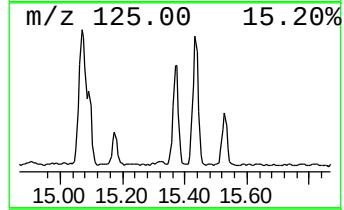
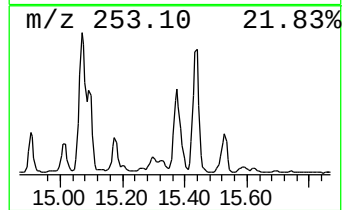
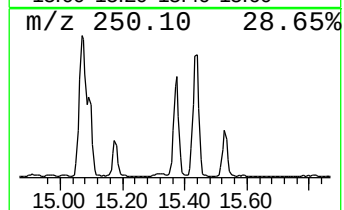
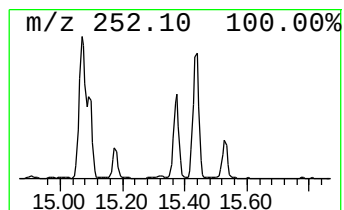
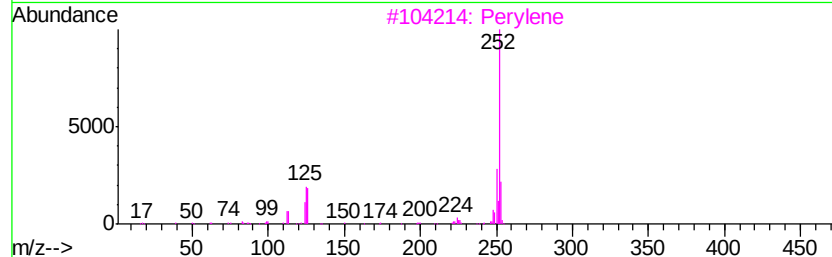
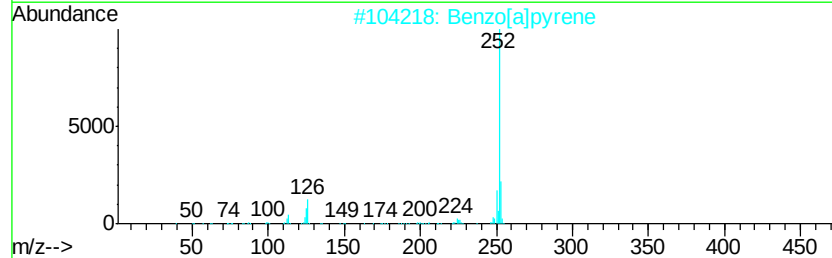
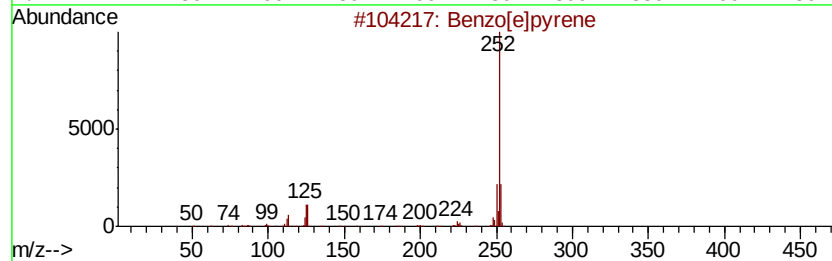
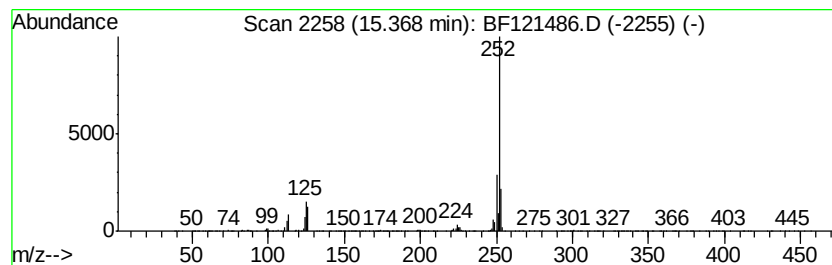
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Perylene Concentration Rank 1

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 15.37 | 15.76 ng | 1498940 | Perylene-d12     | 15.50 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[e]pyrene           | 252 | C20H12  | 000192-97-2 | 99   |
| 2    |    |   | Benzo[a]pyrene           | 252 | C20H12  | 000050-32-8 | 98   |
| 3    |    |   | Perylene                 | 252 | C20H12  | 000198-55-0 | 95   |
| 4    |    |   | Benzo[il]fluoranthene    | 252 | C20H12  | 000205-82-3 | 95   |
| 5    |    |   | Benz[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF083120\  
 Data File : BF121486.D  
 Acq On : 31 Aug 2020 19:32  
 Operator : JU/CG  
 Sample : L3847-15  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 LINDEN-JOP-BH-10-B

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 |
| Butane, 2-methoxy... | 2.15  | 15.5    | ng    | 1431190  | 1                     | 6.86  | 1842650 | 20.0 |
| 9H-Fluorene, 1-me... | 10.99 | 2.0     | ng    | 253652   | 4                     | 11.39 | 2473380 | 20.0 |
| Dibenzothiophene     | 11.28 | 2.8     | ng    | 347729   | 4                     | 11.39 | 2473380 | 20.0 |
| 5H-Dibenzo[a,d]cy... | 11.89 | 4.9     | ng    | 603531   | 4                     | 11.39 | 2473380 | 20.0 |
| Phenanthrene, 1-m... | 11.92 | 8.4     | ng    | 1036930  | 4                     | 11.39 | 2473380 | 20.0 |
| 1H-Cyclopropa[l]p... | 11.96 | 2.8     | ng    | 342659   | 4                     | 11.39 | 2473380 | 20.0 |
| 4H-Cyclopenta[def... | 12.00 | 8.6     | ng    | 1068200  | 4                     | 11.39 | 2473380 | 20.0 |
| Phenanthrene, 2-m... | 12.02 | 2.7     | ng    | 336030   | 4                     | 11.39 | 2473380 | 20.0 |
| Naphthalene, 2-ph... | 12.18 | 3.1     | ng    | 385143   | 4                     | 11.39 | 2473380 | 20.0 |
| Phenanthrene, 2,5... | 12.43 | 2.8     | ng    | 349563   | 4                     | 11.39 | 2473380 | 20.0 |
| unknown12.47         | 12.47 | 2.5     | ng    | 305312   | 4                     | 11.39 | 2473380 | 20.0 |
| Cyclopenta(def)ph... | 12.52 | 2.3     | ng    | 288097   | 4                     | 11.39 | 2473380 | 20.0 |
| Pyrene, 1-methyl-    | 13.04 | 2.3     | ng    | 558462   | 5                     | 14.03 | 4825890 | 20.0 |
| 11H-Benzo[b]fluorene | 13.16 | 4.2     | ng    | 1002850  | 5                     | 14.03 | 4825890 | 20.0 |
| Fluoranthene, 2-m... | 13.26 | 2.0     | ng    | 490474   | 5                     | 14.03 | 4825890 | 20.0 |
| Benzo[c]phenanthrene | 13.80 | 2.1     | ng    | 498115   | 5                     | 14.03 | 4825890 | 20.0 |
| Cyclopenta[cd]pyrene | 13.83 | 2.2     | ng    | 531560   | 5                     | 14.03 | 4825890 | 20.0 |
| 11H-Benzo[a]fluor... | 13.87 | 3.6     | ng    | 879208   | 5                     | 14.03 | 4825890 | 20.0 |
| Benzo[e]pyrene       | 15.17 | 5.8     | ng    | 547209   | 6                     | 15.50 | 1901690 | 20.0 |
| Perylene             | 15.37 | 15.8    | ng    | 1498940  | 6                     | 15.50 | 1901690 | 20.0 |