

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
 Data File : BF130312.D  
 Acq On : 17 Sep 2022 02:41  
 Operator : CG\JU  
 Sample : N4665-02 10X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 COMP-2

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-BF091422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130312.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 6.493       | 746           | 749         | 756          | rVB3     | 1177006        | 1267434       | 6.24%           | 0.406%        |
| 2         | 6.851       | 806           | 810         | 813          | rVB      | 3134338        | 2725514       | 13.43%          | 0.872%        |
| 3         | 6.928       | 820           | 823         | 829          | rBV      | 1522893        | 1510204       | 7.44%           | 0.483%        |
| 4         | 7.704       | 948           | 955         | 958          | rBV3     | 1533729        | 2218515       | 10.93%          | 0.710%        |
| 5         | 7.740       | 958           | 961         | 966          | rVB      | 1631358        | 2138350       | 10.53%          | 0.684%        |
| 6         | 8.004       | 992           | 1006        | 1008         | rBV      | 10453257       | 20300220      | 100.00%         | 6.497%        |
| 7         | 8.034       | 1008          | 1011        | 1016         | rVB      | 900499         | 684008        | 3.37%           | 0.219%        |
| 8         | 8.622       | 1102          | 1111        | 1114         | rVV3     | 409954         | 744764        | 3.67%           | 0.238%        |
| 9         | 8.686       | 1114          | 1122        | 1125         | rVB      | 10118620       | 16392152      | 80.75%          | 5.246%        |
| 10        | 8.786       | 1131          | 1139        | 1141         | rBV      | 8338701        | 11417939      | 56.25%          | 3.654%        |
| 11        | 9.128       | 1190          | 1197        | 1202         | rVB      | 2189911        | 2195188       | 10.81%          | 0.703%        |
| 12        | 9.222       | 1210          | 1213        | 1219         | rVB      | 1787349        | 2318264       | 11.42%          | 0.742%        |
| 13        | 9.304       | 1219          | 1227        | 1229         | rBV      | 4346047        | 6366185       | 31.36%          | 2.038%        |
| 14        | 9.375       | 1233          | 1239        | 1241         | rVV      | 5480232        | 7074469       | 34.85%          | 2.264%        |
| 15        | 9.404       | 1241          | 1244        | 1247         | rVV      | 4639788        | 4197930       | 20.68%          | 1.344%        |
| 16        | 9.486       | 1253          | 1258        | 1266         | rVB      | 3210351        | 3835874       | 18.90%          | 1.228%        |
| 17        | 9.569       | 1266          | 1272        | 1275         | rBV      | 5405811        | 4559770       | 22.46%          | 1.459%        |
| 18        | 9.686       | 1289          | 1292        | 1297         | rBV2     | 1905611        | 2088419       | 10.29%          | 0.668%        |
| 19        | 9.739       | 1297          | 1301        | 1304         | rBV2     | 2240025        | 2264466       | 11.15%          | 0.725%        |
| 20        | 9.810       | 1309          | 1313        | 1317         | rBV2     | 1036263        | 1297714       | 6.39%           | 0.415%        |
| 21        | 9.904       | 1323          | 1329        | 1333         | rVV2     | 2424008        | 2669567       | 13.15%          | 0.854%        |
| 22        | 9.945       | 1333          | 1336        | 1339         | rVV      | 1922784        | 1605032       | 7.91%           | 0.514%        |
| 23        | 10.022      | 1344          | 1349        | 1351         | rBV2     | 1748598        | 2084832       | 10.27%          | 0.667%        |
| 24        | 10.051      | 1351          | 1354        | 1359         | rVB      | 1572819        | 1349559       | 6.65%           | 0.432%        |
| 25        | 10.110      | 1359          | 1364        | 1366         | rBV3     | 1197627        | 1928741       | 9.50%           | 0.617%        |
| 26        | 10.128      | 1366          | 1367        | 1370         | rVV      | 1105920        | 796938        | 3.93%           | 0.255%        |
| 27        | 10.157      | 1370          | 1372        | 1375         | rVB      | 1016086        | 734526        | 3.62%           | 0.235%        |
| 28        | 10.210      | 1375          | 1381        | 1385         | rBV4     | 856608         | 1948026       | 9.60%           | 0.623%        |
| 29        | 10.263      | 1385          | 1390        | 1394         | rVB      | 5848953        | 6840740       | 33.70%          | 2.189%        |
| 30        | 10.316      | 1394          | 1399        | 1401         | rBV3     | 972100         | 1267634       | 6.24%           | 0.406%        |
| 31        | 10.363      | 1404          | 1407        | 1411         | rVV2     | 1163932        | 1234743       | 6.08%           | 0.395%        |
| 32        | 10.404      | 1411          | 1414        | 1417         | rVV2     | 1071880        | 1304775       | 6.43%           | 0.418%        |
| 33        | 10.475      | 1422          | 1426        | 1432         | rVB2     | 2060182        | 2735909       | 13.48%          | 0.876%        |
| 34        | 10.610      | 1447          | 1449        | 1455         | rVB2     | 901861         | 952637        | 4.69%           | 0.305%        |
| 35        | 10.686      | 1457          | 1462        | 1464         | rBV2     | 518017         | 671524        | 3.31%           | 0.215%        |
| 36        | 10.798      | 1475          | 1481        | 1484         | rBV      | 2575992        | 3410947       | 16.80%          | 1.092%        |

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 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 COMP-2

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-BF091422.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |          |          |        |        |
|----|--------|------|------|------|------|----------|----------|--------|--------|
| 37 | 10.828 | 1484 | 1486 | 1489 | rVV  | 2096896  | 1681591  | 8.28%  | 0.538% |
| 38 | 10.880 | 1492 | 1495 | 1498 | rVV  | 1679881  | 1519680  | 7.49%  | 0.486% |
| 39 | 10.945 | 1504 | 1506 | 1509 | rVV3 | 668960   | 731431   | 3.60%  | 0.234% |
| 40 | 10.998 | 1512 | 1515 | 1519 | rVB2 | 1141031  | 1121357  | 5.52%  | 0.359% |
| 41 | 11.086 | 1525 | 1530 | 1533 | rVB2 | 2911932  | 2879950  | 14.19% | 0.922% |
| 42 | 11.239 | 1544 | 1556 | 1559 | rBV  | 10684492 | 20063047 | 98.83% | 6.421% |
| 43 | 11.275 | 1559 | 1562 | 1564 | rVV  | 4115724  | 3817576  | 18.81% | 1.222% |
| 44 | 11.298 | 1564 | 1566 | 1569 | rVV2 | 1073132  | 1135471  | 5.59%  | 0.363% |
| 45 | 11.522 | 1600 | 1604 | 1609 | rVB  | 1986053  | 1874262  | 9.23%  | 0.600% |
| 46 | 11.604 | 1614 | 1618 | 1621 | rVB  | 1770710  | 1928811  | 9.50%  | 0.617% |
| 47 | 11.704 | 1627 | 1635 | 1637 | rBV  | 5418358  | 7800875  | 38.43% | 2.497% |
| 48 | 11.733 | 1637 | 1640 | 1643 | rVB  | 7113350  | 7662192  | 37.74% | 2.452% |
| 49 | 11.775 | 1643 | 1647 | 1650 | rBV  | 2164206  | 2553944  | 12.58% | 0.817% |
| 50 | 11.816 | 1650 | 1654 | 1656 | rVV  | 7111003  | 8063012  | 39.72% | 2.581% |
| 51 | 11.839 | 1656 | 1658 | 1661 | rVB  | 5190257  | 4699894  | 23.15% | 1.504% |
| 52 | 11.986 | 1675 | 1683 | 1686 | rBV  | 2349444  | 3399428  | 16.75% | 1.088% |
| 53 | 12.016 | 1686 | 1688 | 1693 | rVB2 | 862799   | 1000674  | 4.93%  | 0.320% |
| 54 | 12.133 | 1706 | 1708 | 1711 | rVB  | 1102591  | 885284   | 4.36%  | 0.283% |
| 55 | 12.169 | 1711 | 1714 | 1716 | rBV  | 2161064  | 1836775  | 9.05%  | 0.588% |
| 56 | 12.192 | 1716 | 1718 | 1721 | rVB  | 1634891  | 1310815  | 6.46%  | 0.420% |
| 57 | 12.251 | 1722 | 1728 | 1730 | rBV  | 4158270  | 5754159  | 28.35% | 1.842% |
| 58 | 12.286 | 1730 | 1734 | 1739 | rVB2 | 3171625  | 5309216  | 26.15% | 1.699% |
| 59 | 12.333 | 1739 | 1742 | 1747 | rVB2 | 1446881  | 2382563  | 11.74% | 0.763% |
| 60 | 12.416 | 1751 | 1756 | 1759 | rBV  | 6896170  | 8716473  | 42.94% | 2.790% |
| 61 | 12.451 | 1759 | 1762 | 1764 | rBV  | 843887   | 945866   | 4.66%  | 0.303% |
| 62 | 12.498 | 1767 | 1770 | 1773 | rVB  | 1300752  | 1189924  | 5.86%  | 0.381% |
| 63 | 12.551 | 1775 | 1779 | 1781 | rBV2 | 676269   | 672516   | 3.31%  | 0.215% |
| 64 | 12.657 | 1789 | 1797 | 1799 | rBV  | 9081363  | 14855932 | 73.18% | 4.755% |
| 65 | 12.686 | 1799 | 1802 | 1806 | rVB2 | 1471970  | 1679581  | 8.27%  | 0.538% |
| 66 | 12.745 | 1806 | 1812 | 1814 | rBV4 | 814338   | 1390120  | 6.85%  | 0.445% |
| 67 | 12.786 | 1817 | 1819 | 1824 | rVB3 | 932247   | 1114157  | 5.49%  | 0.357% |
| 68 | 12.851 | 1825 | 1830 | 1836 | rBV2 | 2137520  | 3238808  | 15.95% | 1.037% |
| 69 | 12.904 | 1836 | 1839 | 1841 | rBV2 | 614889   | 651497   | 3.21%  | 0.209% |
| 70 | 12.939 | 1841 | 1845 | 1846 | rBV2 | 1377389  | 1539029  | 7.58%  | 0.493% |
| 71 | 12.969 | 1846 | 1850 | 1852 | rVB  | 3655619  | 4068611  | 20.04% | 1.302% |
| 72 | 13.033 | 1852 | 1861 | 1863 | rBV  | 2814631  | 4090074  | 20.15% | 1.309% |
| 73 | 13.069 | 1863 | 1867 | 1870 | rBV2 | 3074792  | 3231749  | 15.92% | 1.034% |
| 74 | 13.127 | 1874 | 1877 | 1879 | rBV2 | 761536   | 849480   | 4.18%  | 0.272% |
| 75 | 13.163 | 1879 | 1883 | 1885 | rBV  | 2170561  | 1928641  | 9.50%  | 0.617% |
| 76 | 13.192 | 1885 | 1888 | 1891 | rVB  | 3188203  | 2913667  | 14.35% | 0.933% |

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 Sample : N4665-02 10X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 COMP-2

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-BF091422.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 77  | 13.363 | 1912 | 1917 | 1918 | rBV2 | 742147  | 908812  | 4.48%  | 0.291% |
| 78  | 13.439 | 1926 | 1930 | 1933 | rBV2 | 1027824 | 1625760 | 8.01%  | 0.520% |
| 79  | 13.580 | 1947 | 1954 | 1956 | rVV2 | 1413000 | 2429980 | 11.97% | 0.778% |
| 80  | 13.604 | 1956 | 1958 | 1960 | rVV  | 1483413 | 1314562 | 6.48%  | 0.421% |
| 81  | 13.627 | 1960 | 1962 | 1964 | rVV  | 800582  | 739966  | 3.65%  | 0.237% |
| 82  | 13.821 | 1989 | 1995 | 1998 | rBV  | 4322151 | 6926995 | 34.12% | 2.217% |
| 83  | 13.863 | 1998 | 2002 | 2005 | rVV  | 4581724 | 5797030 | 28.56% | 1.855% |
| 84  | 14.174 | 2052 | 2055 | 2057 | rBV  | 587217  | 695407  | 3.43%  | 0.223% |
| 85  | 14.216 | 2057 | 2062 | 2065 | rVV  | 2537417 | 3073510 | 15.14% | 0.984% |
| 86  | 14.245 | 2065 | 2067 | 2070 | rVB  | 1664206 | 1268054 | 6.25%  | 0.406% |
| 87  | 14.304 | 2075 | 2077 | 2080 | rVV2 | 1195485 | 1285583 | 6.33%  | 0.411% |
| 88  | 14.339 | 2080 | 2083 | 2085 | rVV  | 1354215 | 1585639 | 7.81%  | 0.507% |
| 89  | 14.357 | 2085 | 2086 | 2090 | rVV  | 1278771 | 956412  | 4.71%  | 0.306% |
| 90  | 14.404 | 2090 | 2094 | 2097 | rVV2 | 564857  | 657188  | 3.24%  | 0.210% |
| 91  | 14.533 | 2113 | 2116 | 2122 | rVB3 | 431942  | 850532  | 4.19%  | 0.272% |
| 92  | 14.616 | 2127 | 2130 | 2133 | rVB2 | 616304  | 674431  | 3.32%  | 0.216% |
| 93  | 14.733 | 2148 | 2150 | 2161 | rVB3 | 507951  | 854441  | 4.21%  | 0.273% |
| 94  | 14.833 | 2161 | 2167 | 2170 | rBV  | 1667426 | 3021378 | 14.88% | 0.967% |
| 95  | 14.927 | 2179 | 2183 | 2186 | rVB  | 610678  | 671608  | 3.31%  | 0.215% |
| 96  | 15.110 | 2209 | 2214 | 2218 | rVB  | 1309424 | 1548422 | 7.63%  | 0.496% |
| 97  | 15.174 | 2218 | 2225 | 2228 | rBV  | 2253968 | 2913476 | 14.35% | 0.932% |
| 98  | 15.245 | 2234 | 2237 | 2243 | rVB2 | 478569  | 773682  | 3.81%  | 0.248% |
| 99  | 16.551 | 2453 | 2459 | 2465 | rVB2 | 611225  | 1197007 | 5.90%  | 0.383% |
| 100 | 16.957 | 2521 | 2528 | 2536 | rVB  | 526515  | 1049861 | 5.17%  | 0.336% |

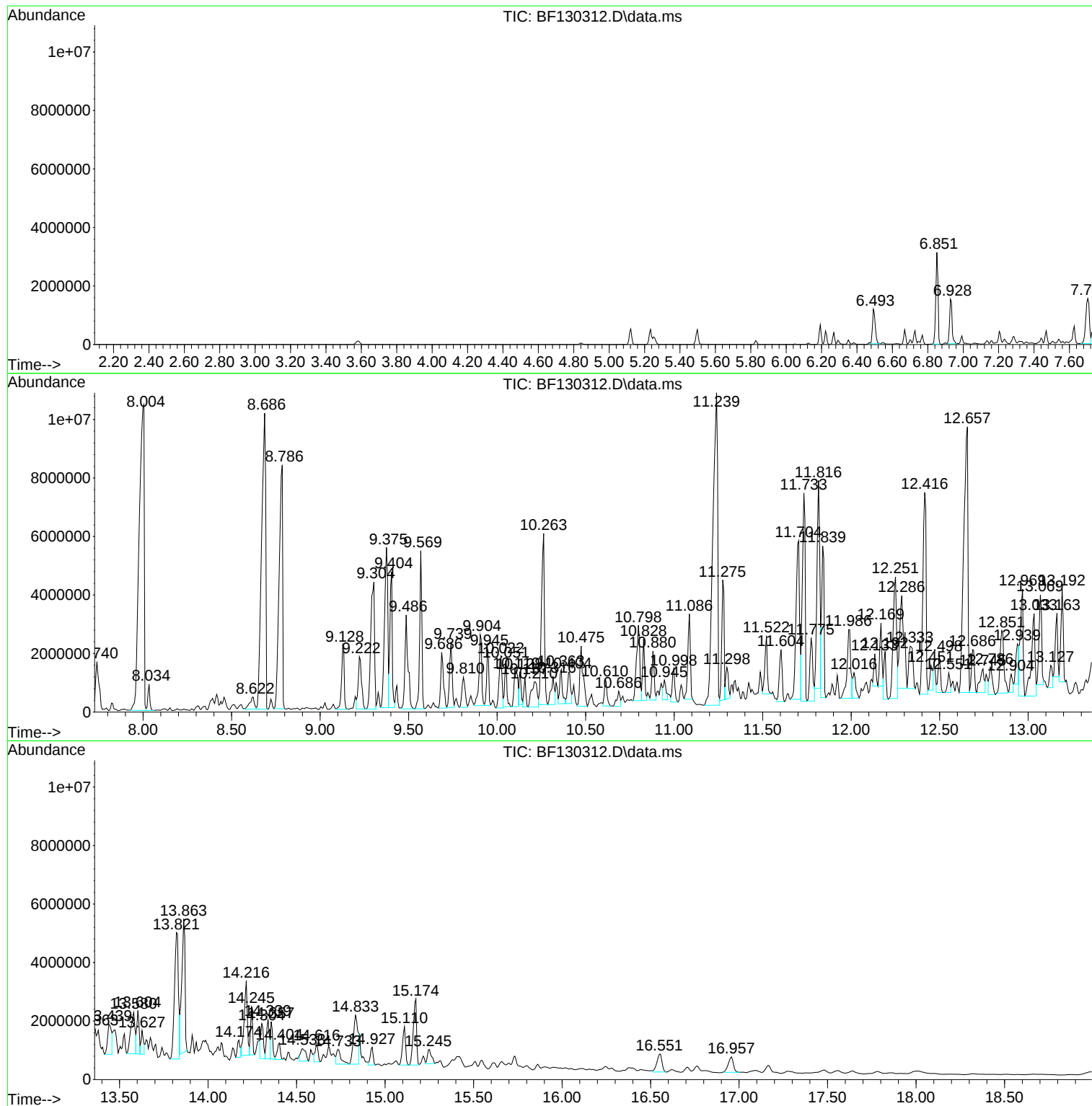
Sum of corrected areas: 312441377

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ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

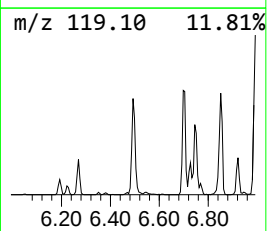
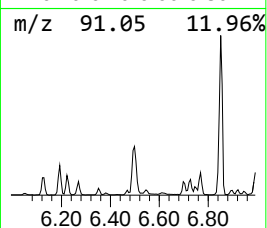
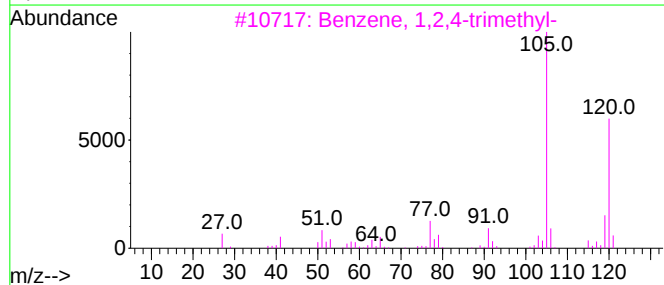
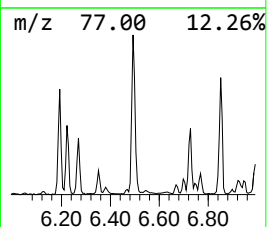
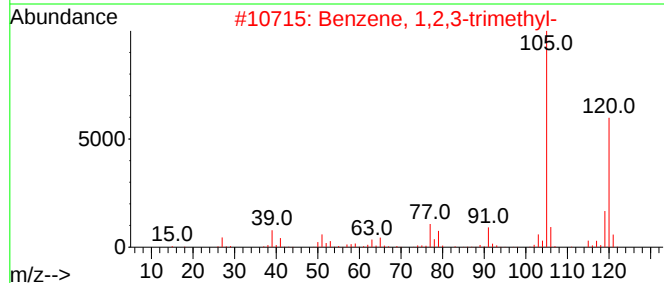
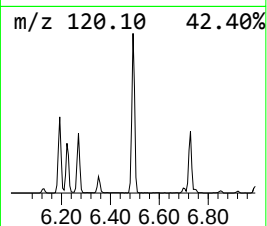
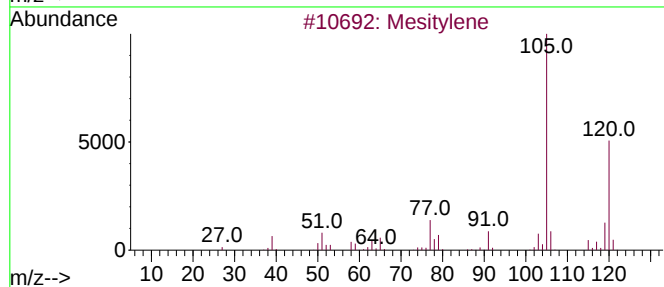
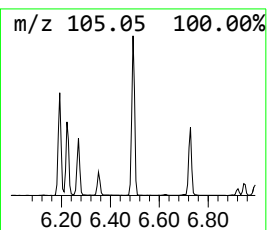
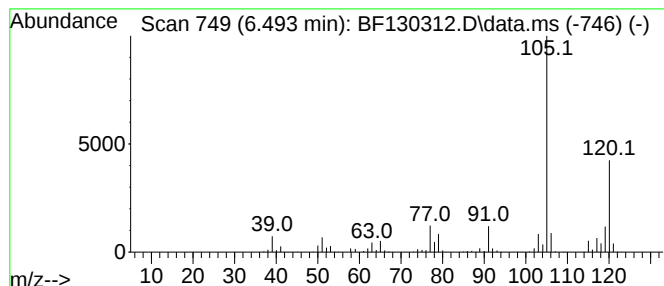
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Mesitylene Concentration Rank 11

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 6.493 | 20.00 ng | 1267430 | 1,4-Dichlorobenzene-d4 | 6.669 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 97   |
| 2    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 93   |
| 3    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 87   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 87   |



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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

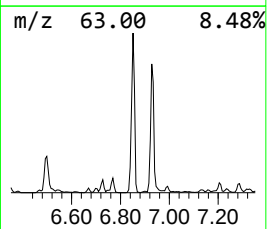
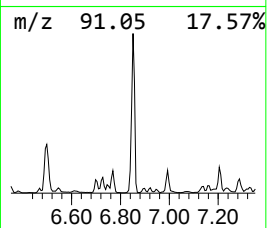
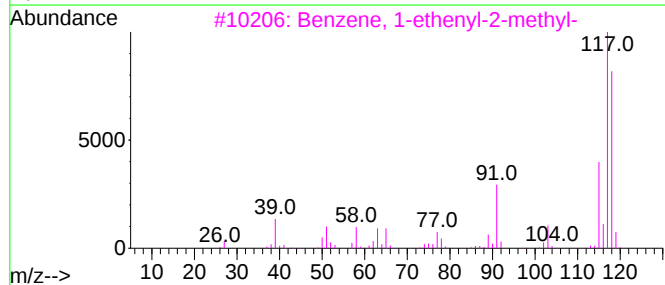
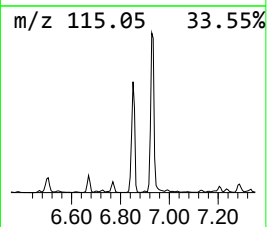
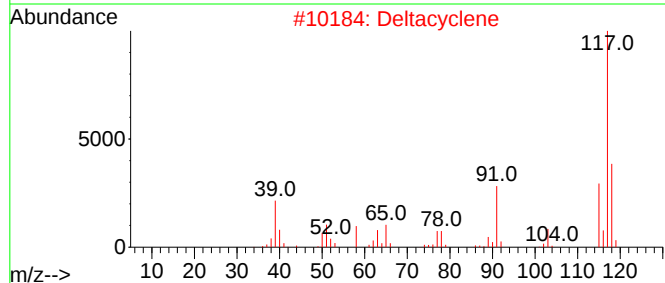
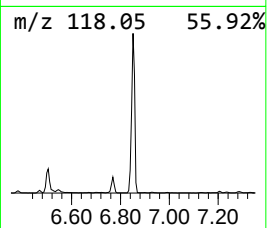
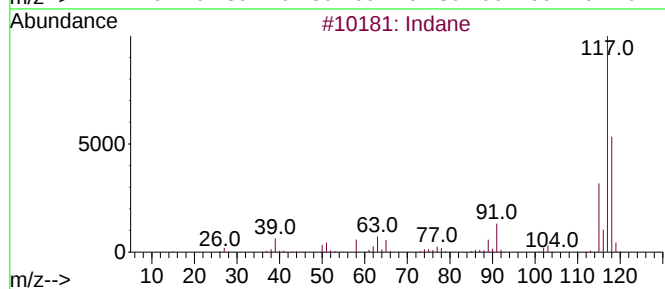
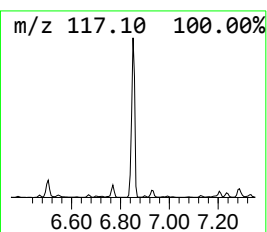
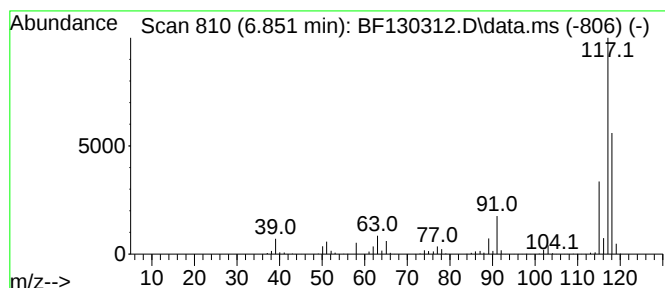
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Indane Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 6.851 | 43.01 ng | 2725510 | 1,4-Dichlorobenzene-d4 | 6.669 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Indane                          | 118 | C9H10   | 000496-11-7 | 91   |
| 2    |    |   | Deltacyclene                    | 118 | C9H10   | 007785-10-6 | 74   |
| 3    |    |   | Benzene, 1-ethenyl-2-methyl-    | 118 | C9H10   | 000611-15-4 | 74   |
| 4    |    |   | Benzene, 2-propenyl-            | 118 | C9H10   | 000300-57-2 | 60   |
| 5    |    |   | Benzeneethanol, .beta.-ethenyl- | 148 | C10H12O | 006052-63-7 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

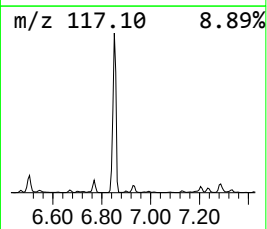
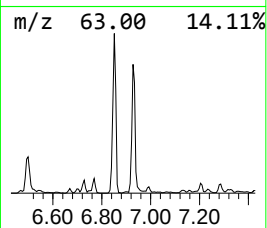
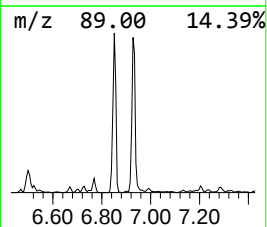
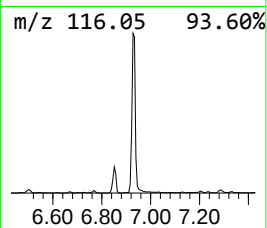
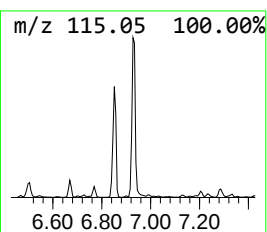
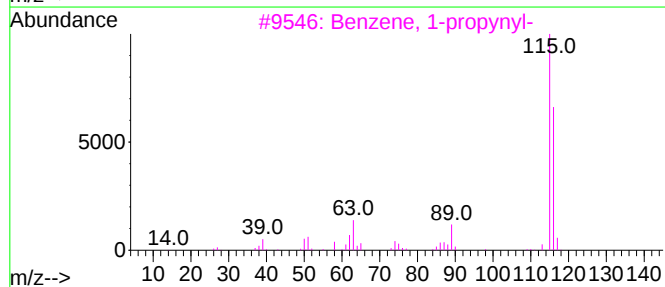
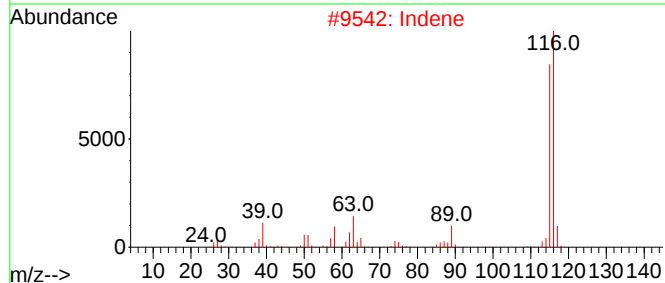
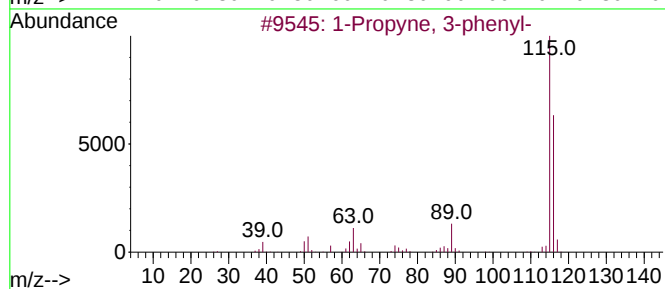
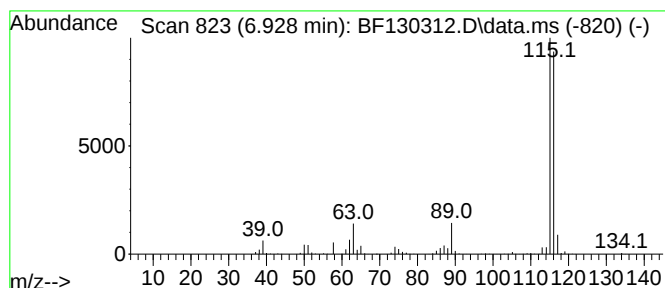
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 1-Propyne, 3-phenyl- Concentration Rank 7

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 6.928 | 23.83 ng | 1510200 | 1,4-Dichlorobenzene-d4 | 6.669 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Propyne, 3-phenyl-      | 116 | C9H8    | 010147-11-2 | 94   |
| 2    |    |   | Indene                    | 116 | C9H8    | 000095-13-6 | 94   |
| 3    |    |   | Benzene, 1-propynyl-      | 116 | C9H8    | 000673-32-5 | 91   |
| 4    |    |   | 3-Methylphenylacetylene   | 116 | C9H8    | 000766-82-5 | 91   |
| 5    |    |   | Benzene, 1,2-propadienyl- | 116 | C9H8    | 002327-99-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

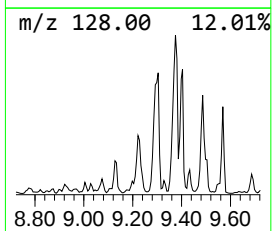
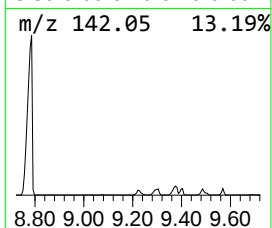
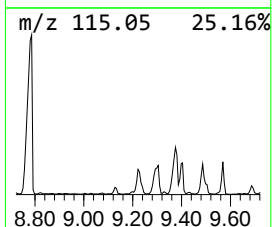
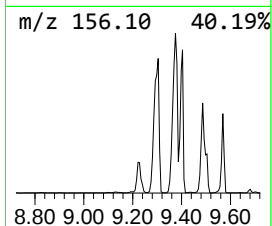
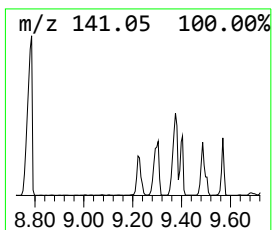
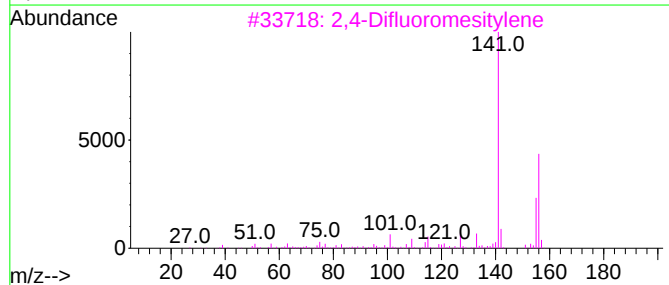
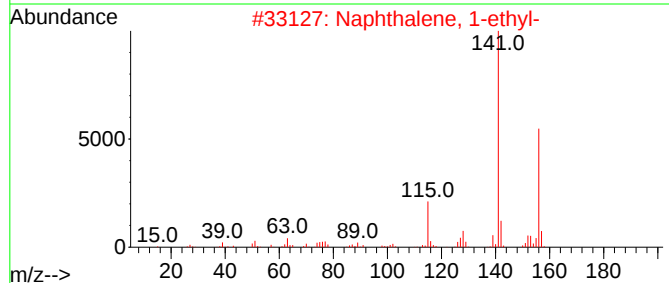
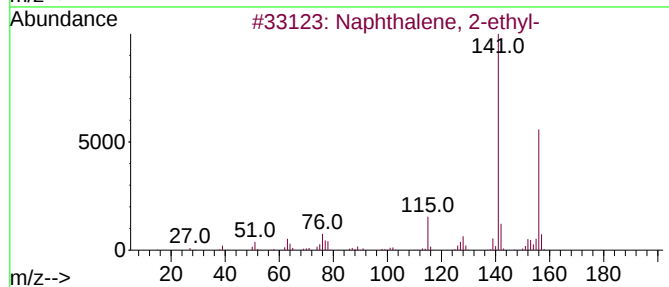
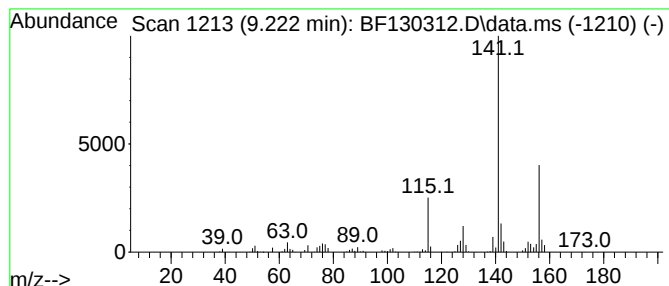
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 Naphthalene, 2-ethyl- Concentration Rank 8

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 9.222 | 22.20 ng | 2318260 | Acenaphthene-d10 | 9.704 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2-ethyl-               | 156 | C12H12  | 000939-27-5 | 96   |
| 2    |    |   | Naphthalene, 1-ethyl-               | 156 | C12H12  | 001127-76-0 | 95   |
| 3    |    |   | 2,4-Difluoromesitylene              | 156 | C9H10F2 | 000392-61-0 | 59   |
| 4    |    |   | Benzenemethanol, 4-chloro-.alpha... | 156 | C8H9ClO | 003391-10-4 | 53   |
| 5    |    |   | Methyl phenylphosphinic acid        | 156 | C7H9O2P | 004271-13-0 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

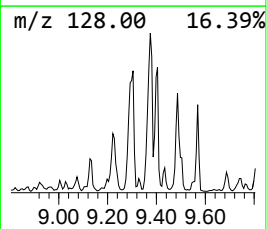
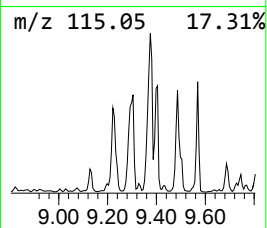
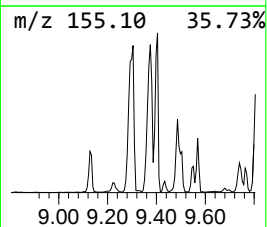
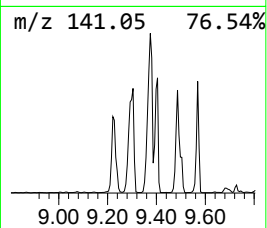
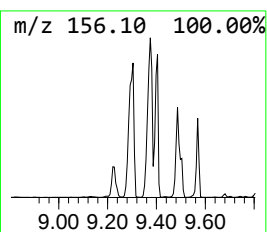
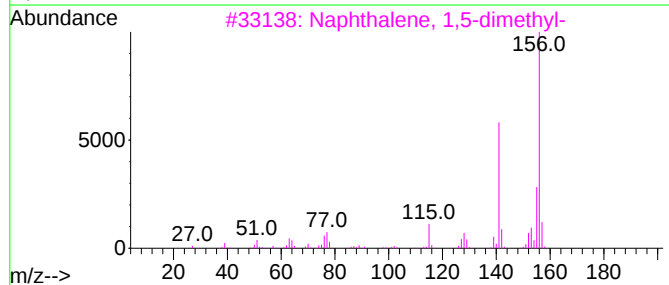
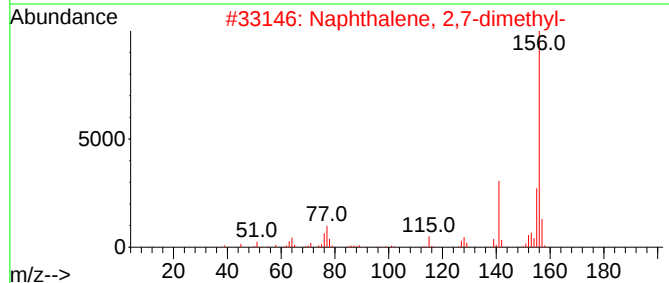
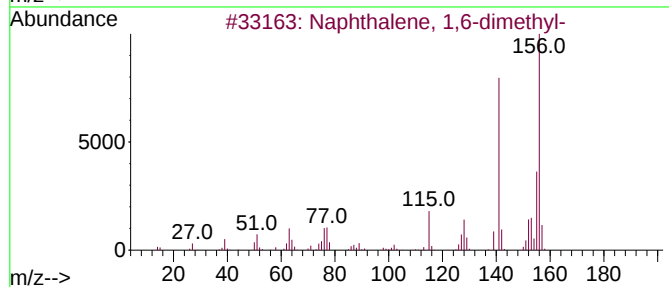
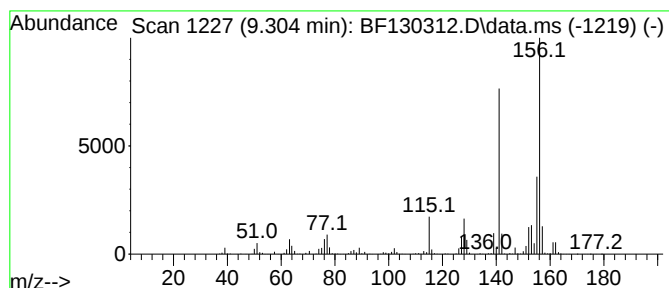
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Naphthalene, 1,6-dimethyl- Concentration Rank 2

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 9.304 | 60.97 ng | 6366190 | Acenaphthene-d10 | 9.704 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 2    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |
| 3    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 96   |
| 4    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 5    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

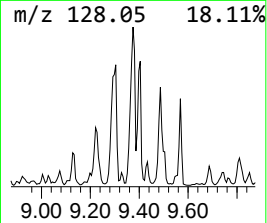
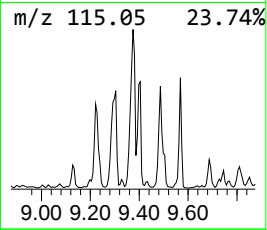
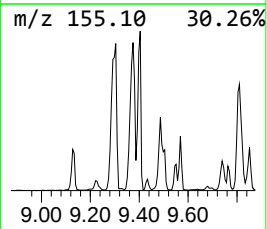
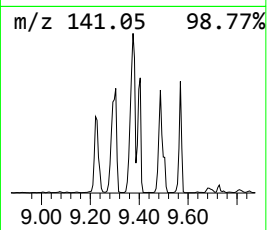
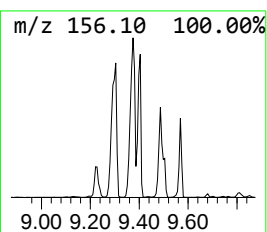
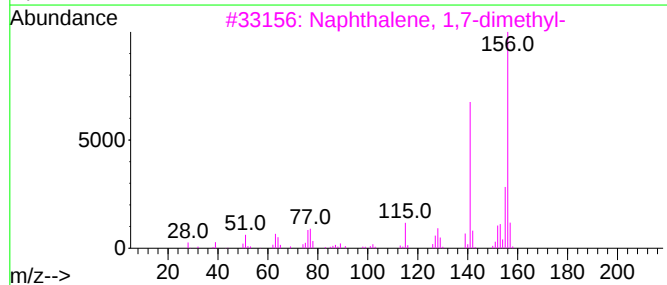
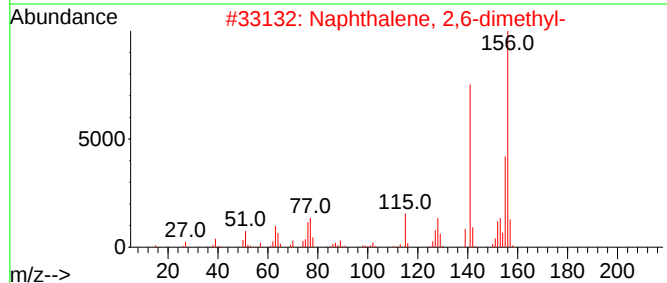
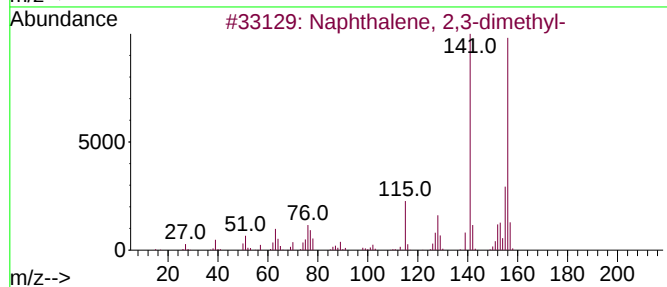
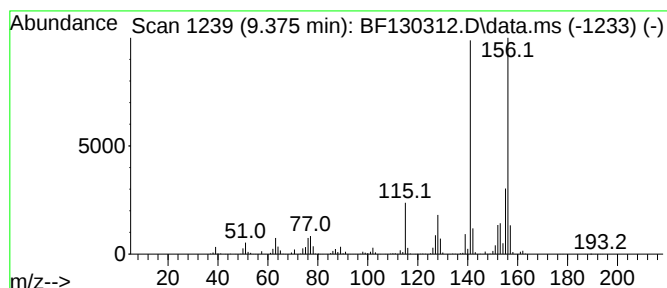
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Naphthalene, 2,3-dimethyl- Concentration Rank 1

| R.T.      | EstConc                    | Area    | Relative to ISTD |             | R.T.  |
|-----------|----------------------------|---------|------------------|-------------|-------|
| 9.375     | 67.75 ng                   | 7074470 | Acenaphthene-d10 |             | 9.704 |
| 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, 1,7-dimethyl- | 156     | C12H12           | 000575-37-1 | 97    |
| 4         | Naphthalene, 1,3-dimethyl- | 156     | C12H12           | 000575-41-7 | 96    |
| 5         | Naphthalene, 1,6-dimethyl- | 156     | C12H12           | 000575-43-9 | 96    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
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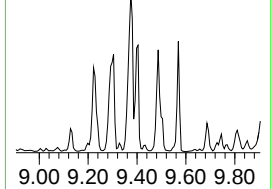
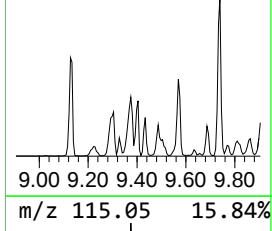
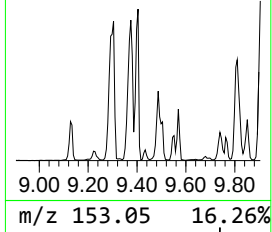
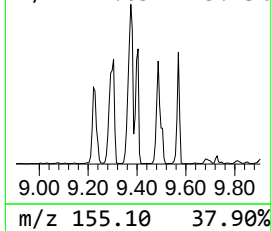
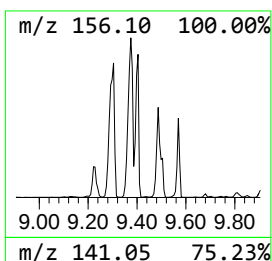
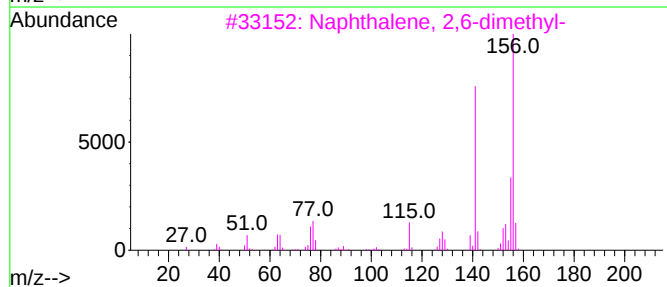
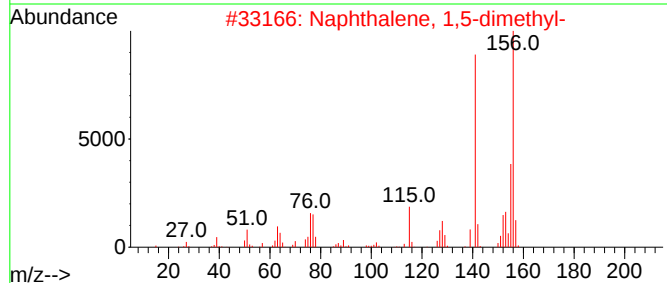
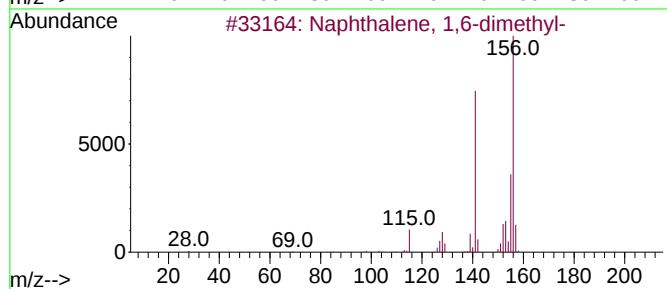
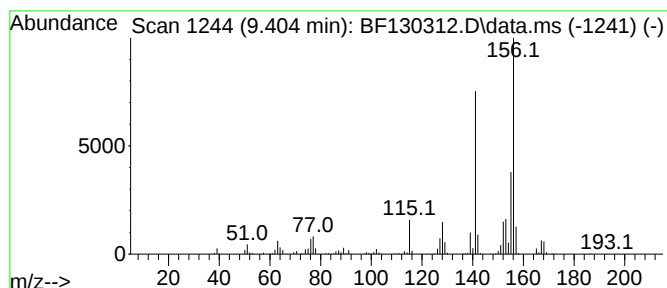
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Peak Number 7 Naphthalene, 1,5-dimethyl- Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 9.404 | 40.20 ng | 4197930 | Acenaphthene-d10 | 9.704 |

| Hit# | of 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 2    |      | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 95   |
| 3    |      | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 95   |
| 4    |      | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 95   |
| 5    |      | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

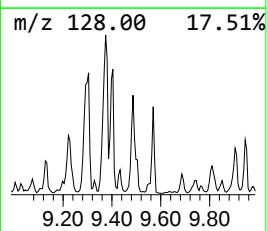
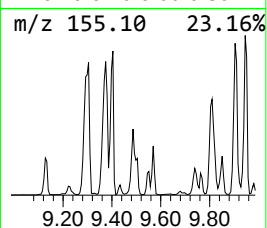
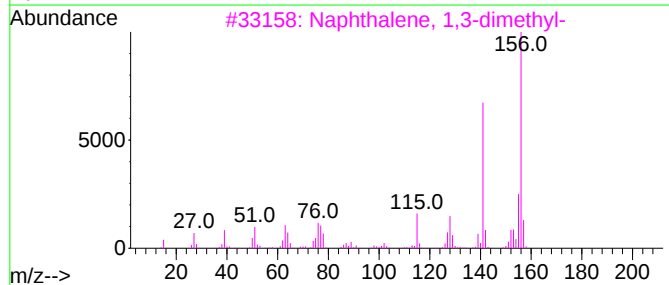
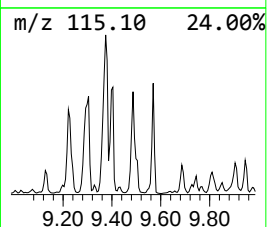
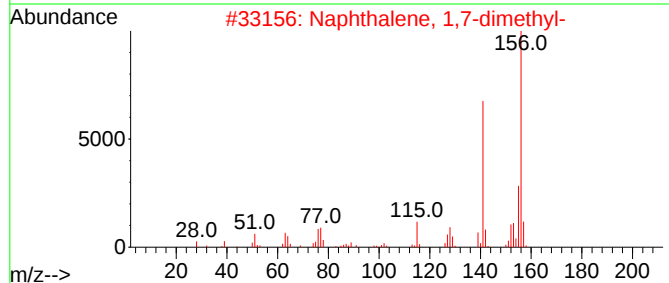
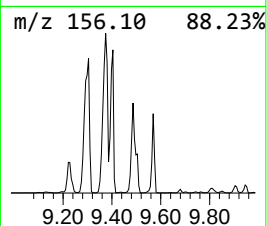
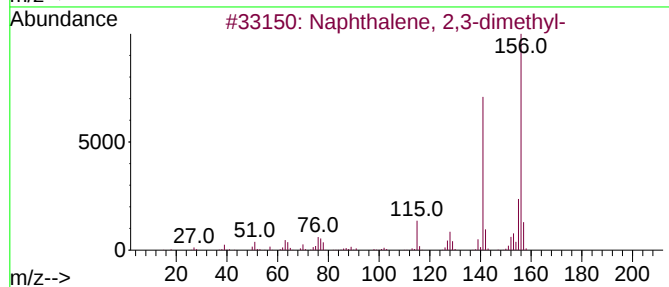
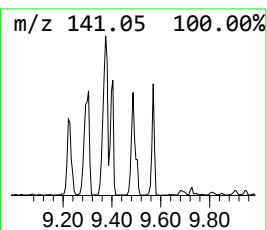
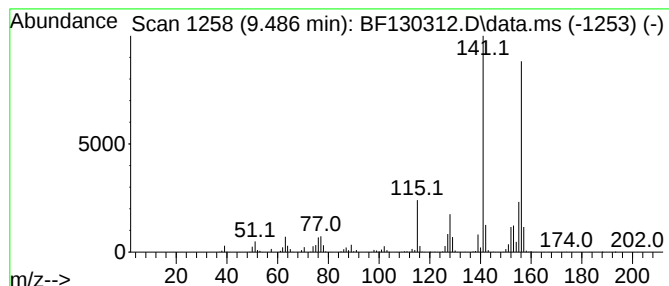
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Naphthalene, 1,7-dimethyl- Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 9.486 | 36.73 ng | 3835870 | Acenaphthene-d10 | 9.704 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 2    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 3    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 97   |
| 4    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 96   |
| 5    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

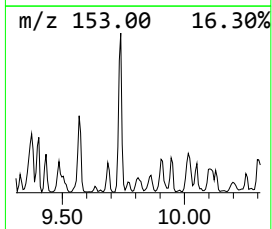
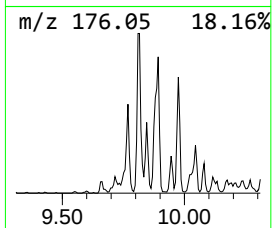
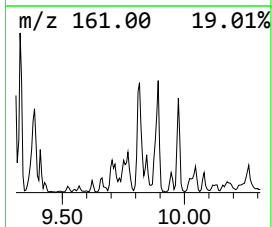
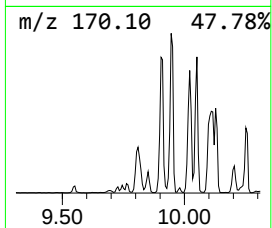
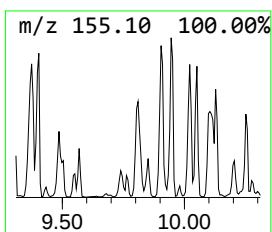
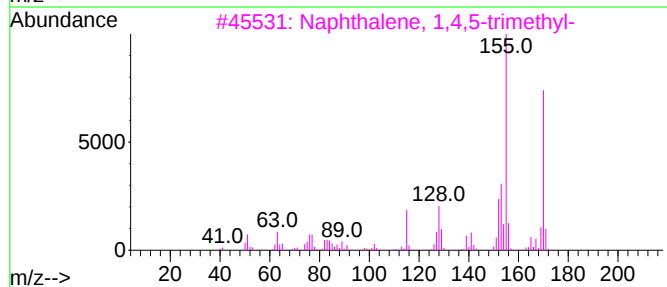
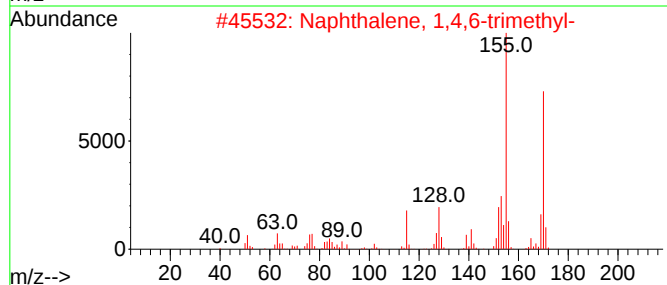
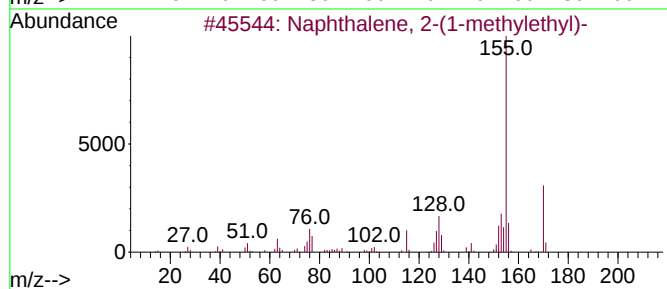
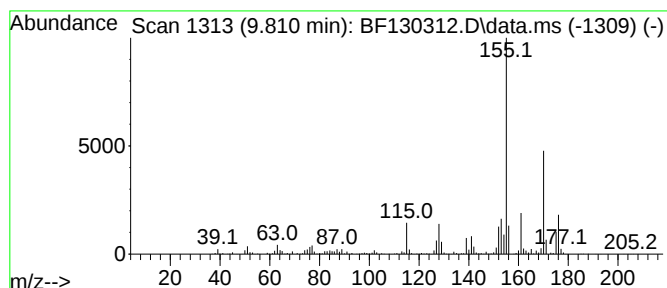
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Naphthalene, 2-(1-methyleth... Concentration Rank 19

| R.T.      | EstConc                         | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|---------------------------------|---------|------------------|---------|-------------|------|
| 9.810     | 12.43 ng                        | 1297710 | Acenaphthene-d10 |         | 9.704       |      |
| Hit# of 5 | Tentative ID                    |         | MW               | MolForm | CAS#        | Qual |
| 1         | Naphthalene, 2-(1-methylethyl)- |         | 170              | C13H14  | 002027-17-0 | 91   |
| 2         | Naphthalene, 1,4,6-trimethyl-   |         | 170              | C13H14  | 002131-42-2 | 91   |
| 3         | Naphthalene, 1,4,5-trimethyl-   |         | 170              | C13H14  | 002131-41-1 | 70   |
| 4         | 2-Naphthyl methyl ketone        |         | 170              | C12H10O | 000093-08-3 | 64   |
| 5         | Ethanone, 1-(1-Naphthalenyl)-   |         | 170              | C12H10O | 000941-98-0 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

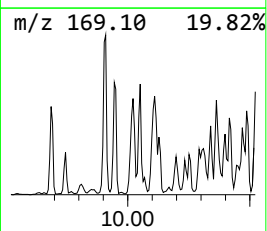
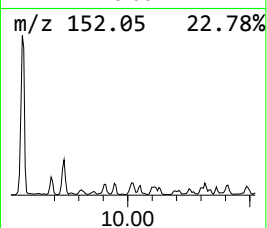
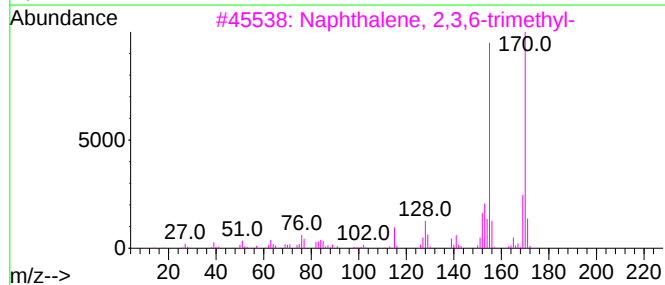
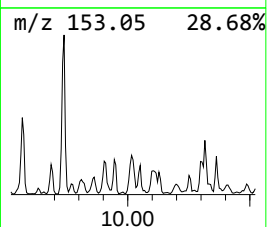
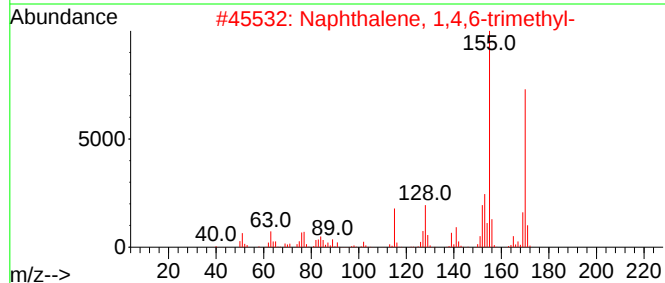
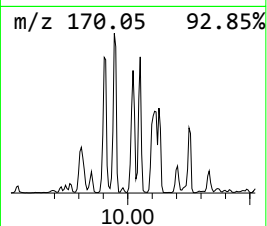
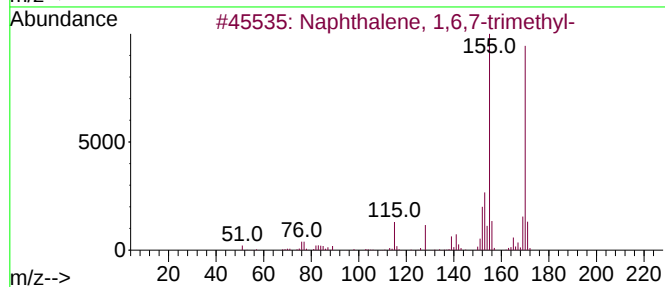
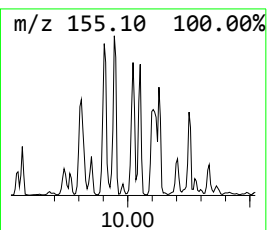
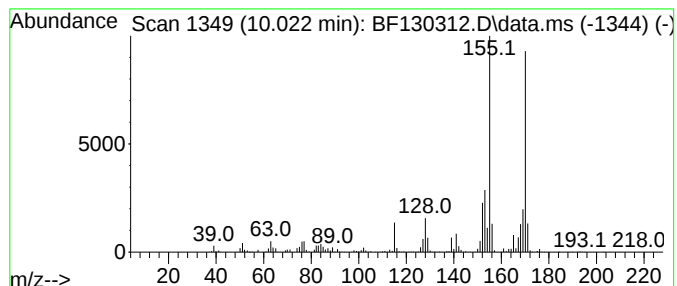
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Naphthalene, 1,6,7-trimethyl- Concentration Rank 12

| R.T.      | EstConc                       | Area    | Relative to ISTD |         |             | R.T.  |
|-----------|-------------------------------|---------|------------------|---------|-------------|-------|
| 10.022    | 19.97 ng                      | 2084830 | Acenaphthene-d10 |         |             | 9.704 |
| Hit# of 5 | Tentative ID                  |         | MW               | MolForm | CAS#        | Qual  |
| 1         | Naphthalene, 1,6,7-trimethyl- |         | 170              | C13H14  | 002245-38-7 | 98    |
| 2         | Naphthalene, 1,4,6-trimethyl- |         | 170              | C13H14  | 002131-42-2 | 97    |
| 3         | Naphthalene, 2,3,6-trimethyl- |         | 170              | C13H14  | 000829-26-5 | 95    |
| 4         | Naphthalene, 1,4,5-trimethyl- |         | 170              | C13H14  | 002131-41-1 | 95    |
| 5         | 4,6,8-Trimethylazulene        |         | 170              | C13H14  | 000941-81-1 | 91    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

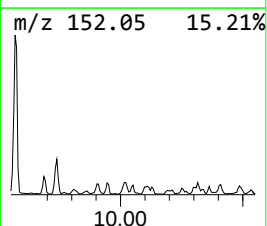
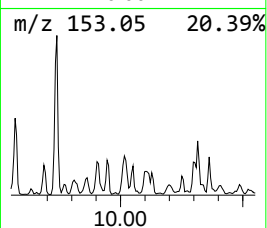
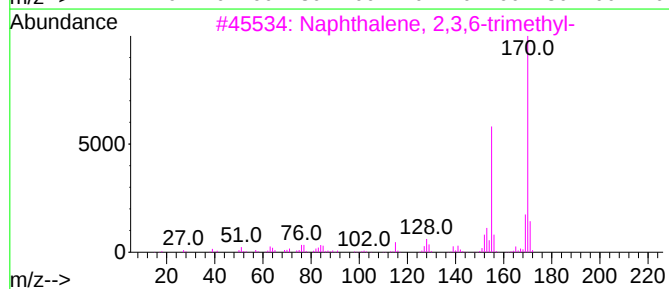
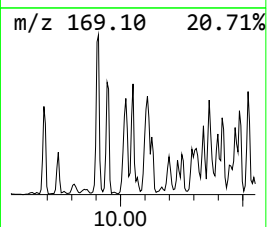
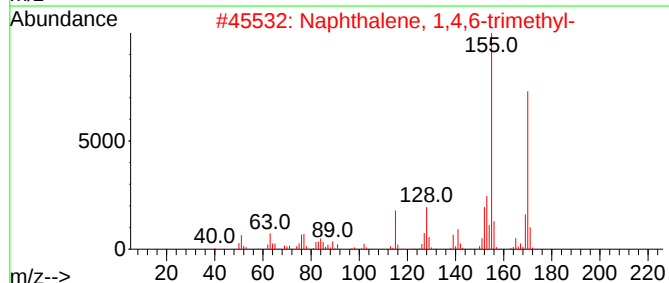
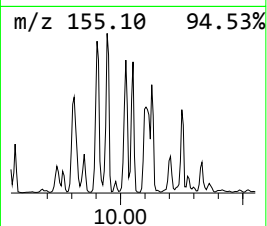
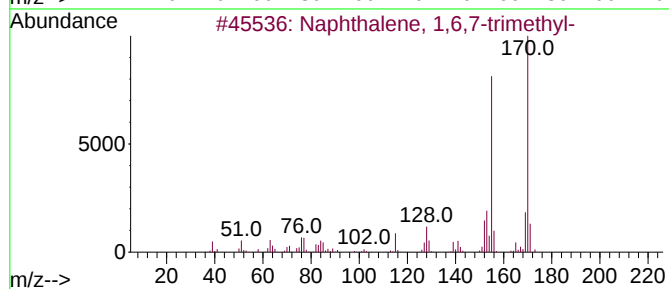
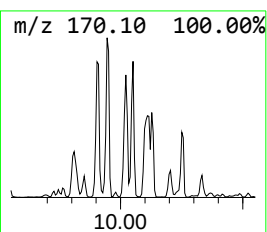
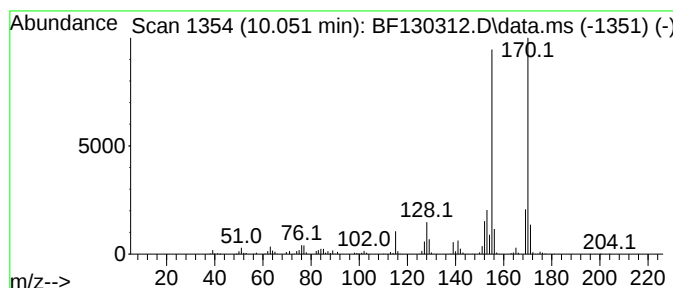
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Naphthalene, 1,4,6-trimethyl- Concentration Rank 17

| R.T.      | EstConc                             | Area    | Relative to ISTD |         | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|---------|--------------|------|
| 10.051    | 12.92 ng                            | 1349560 | Acenaphthene-d10 |         | 9.704        |      |
| Hit# of 5 | Tentative ID                        |         | MW               | MolForm | CAS#         | Qual |
| 1         | Naphthalene, 1,6,7-trimethyl-       |         | 170              | C13H14  | 002245-38-7  | 97   |
| 2         | Naphthalene, 1,4,6-trimethyl-       |         | 170              | C13H14  | 002131-42-2  | 97   |
| 3         | Naphthalene, 2,3,6-trimethyl-       |         | 170              | C13H14  | 000829-26-5  | 96   |
| 4         | Naphthalene, 1,4,5-trimethyl-       |         | 170              | C13H14  | 002131-41-1  | 94   |
| 5         | 1-(2,2-Dimethylcyclopropyl)-2-ph... |         | 170              | C13H14  | 1000294-91-1 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

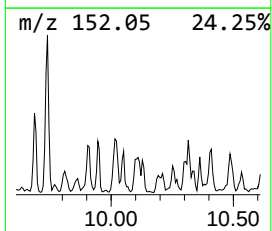
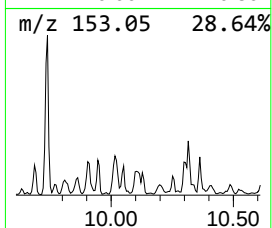
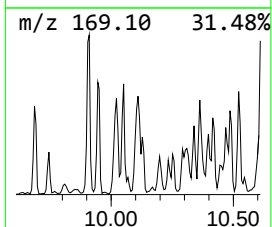
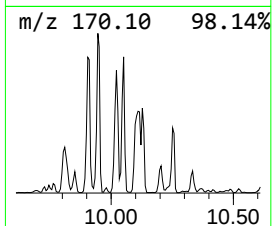
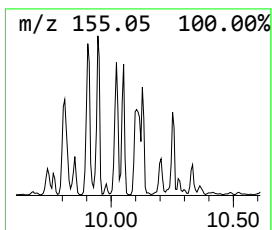
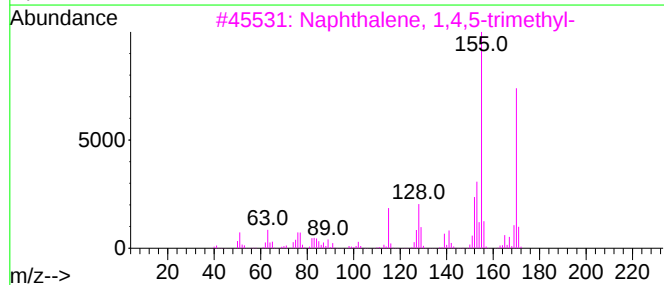
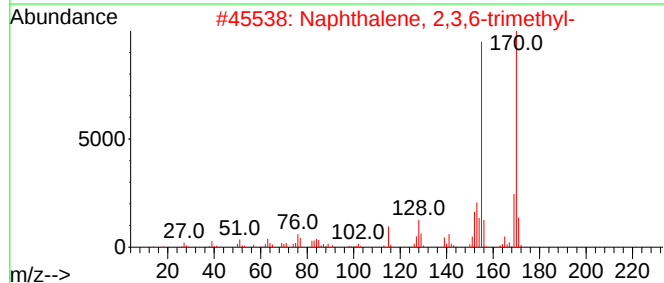
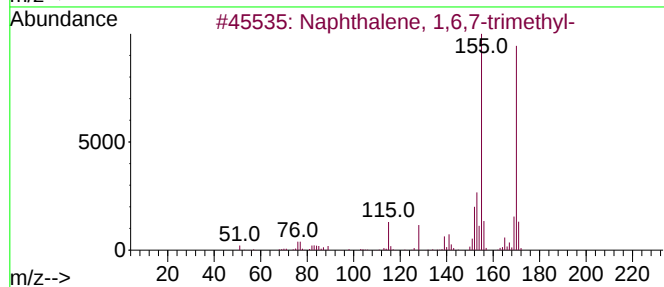
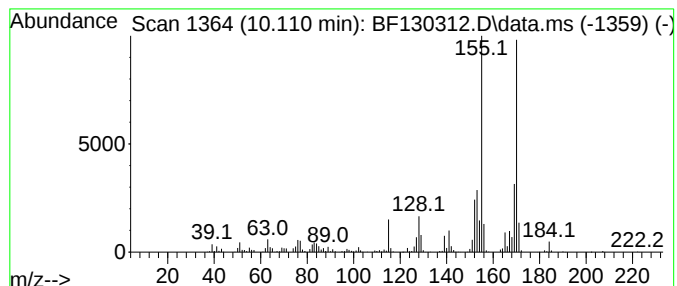
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 Naphthalene, 2,3,6-trimethyl- Concentration Rank 14

| R.T.      | EstConc                       | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|-------------------------------|---------|------------------|---------|-------------|------|
| 10.110    | 18.47 ng                      | 1928740 | Acenaphthene-d10 |         | 9.704       |      |
| Hit# of 5 | Tentative ID                  |         | MW               | MolForm | CAS#        | Qual |
| 1         | Naphthalene, 1,6,7-trimethyl- |         | 170              | C13H14  | 002245-38-7 | 96   |
| 2         | Naphthalene, 2,3,6-trimethyl- |         | 170              | C13H14  | 000829-26-5 | 96   |
| 3         | Naphthalene, 1,4,5-trimethyl- |         | 170              | C13H14  | 002131-41-1 | 96   |
| 4         | Naphthalene, 1,4,6-trimethyl- |         | 170              | C13H14  | 002131-42-2 | 95   |
| 5         | 4,6,8-Trimethylazulene        |         | 170              | C13H14  | 000941-81-1 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
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TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

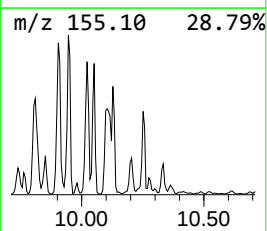
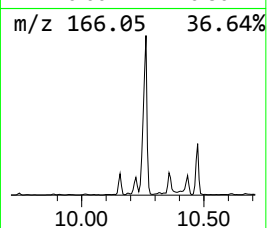
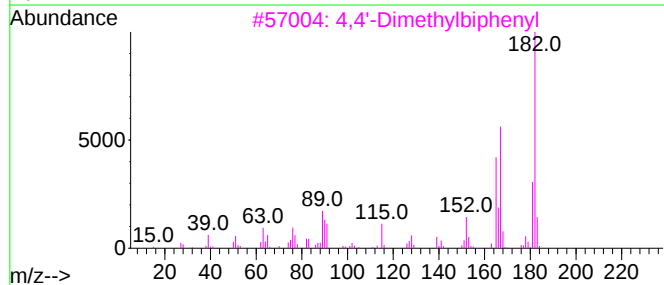
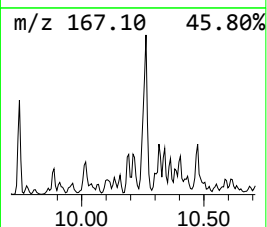
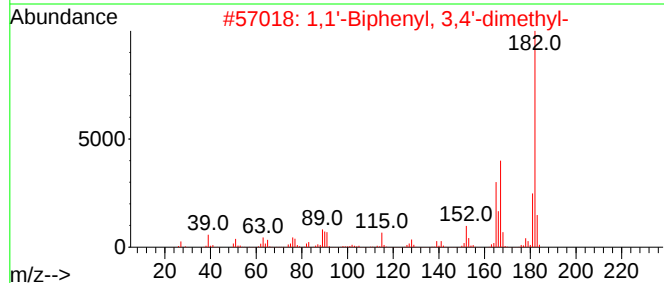
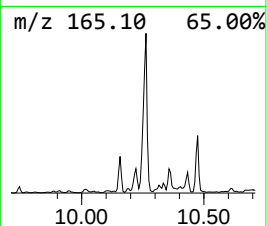
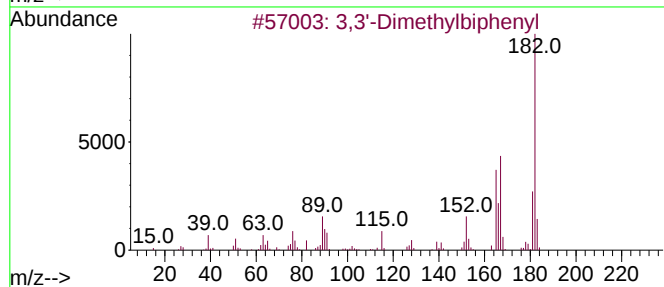
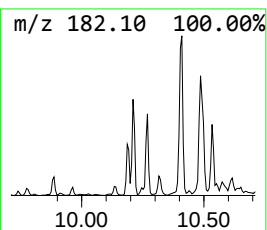
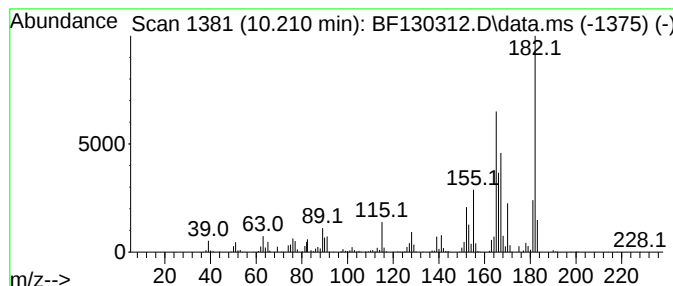
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Peak Number 13 3,3'-Dimethylbiphenyl Concentration Rank 13

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.  |
|--------|----------|---------|------------------|-------|
| 10.210 | 18.66 ng | 1948030 | Acenaphthene-d10 | 9.704 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 3,3'-Dimethylbiphenyl               | 182 | C14H14  | 000612-75-9 | 94   |
| 2    |      | 1,1'-Biphenyl, 3,4'-dimethyl-       | 182 | C14H14  | 007383-90-6 | 70   |
| 3    |      | 4,4'-Dimethylbiphenyl               | 182 | C14H14  | 000613-33-2 | 70   |
| 4    |      | Dehydrochamazulene                  | 182 | C14H14  | 321732-25-6 | 60   |
| 5    |      | Bicyclo[3.2.1]octa-2,6-diene, 2-... | 182 | C14H14  | 101166-08-9 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

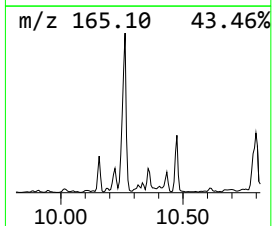
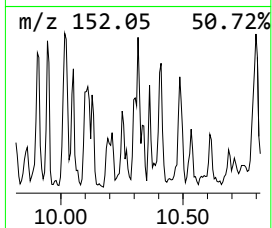
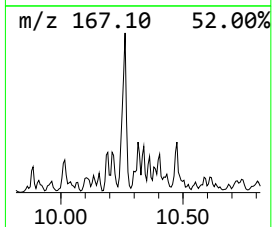
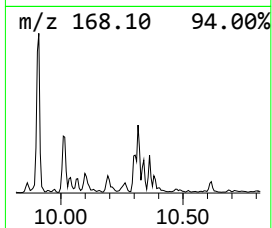
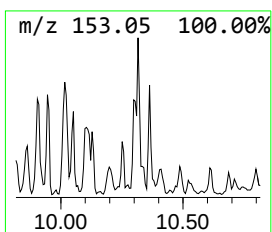
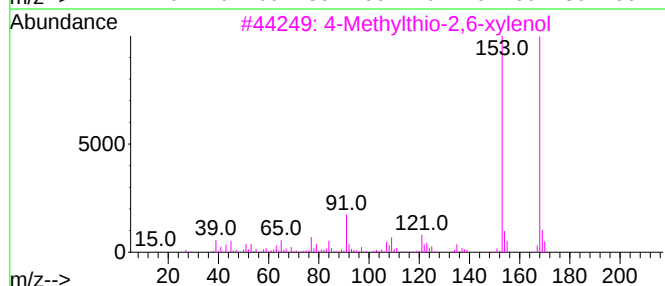
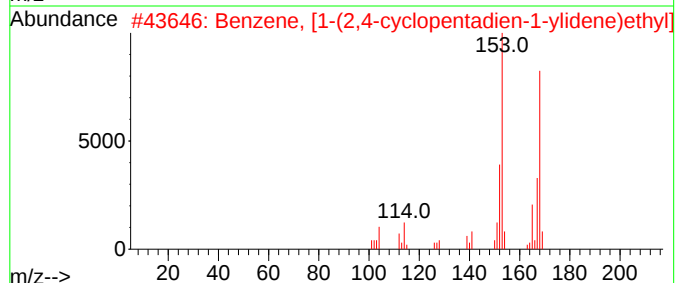
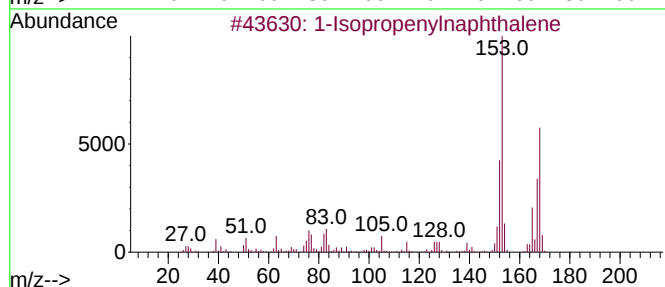
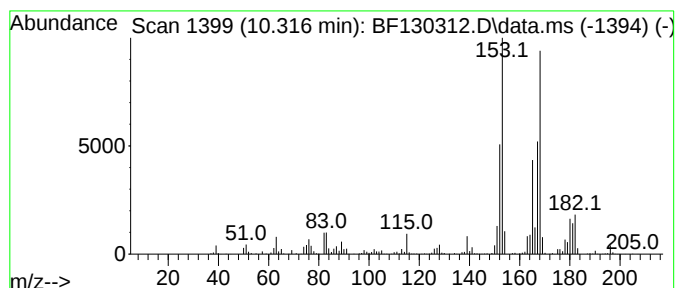
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 1-Isopropenylnaphthalene Concentration Rank 20

| R.T.   | EstConc  | Area                                | Relative to ISTD | R.T.           |
|--------|----------|-------------------------------------|------------------|----------------|
| 10.316 | 12.14 ng | 1267630                             | Acenaphthene-d10 | 9.704          |
| Hit#   | of 5     | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1      |          | 1-Isopropenylnaphthalene            | 168 C13H12       | 001855-47-6 89 |
| 2      |          | Benzene, [1-(2,4-cyclopentadien-... | 168 C13H12       | 002320-32-3 70 |
| 3      |          | 4-Methylthio-2,6-xyleneol           | 168 C9H12OS      | 007379-49-9 43 |
| 4      |          | 1,1'-Biphenyl, 2-methyl-            | 168 C13H12       | 000643-58-3 38 |
| 5      |          | 1,1'-Biphenyl, 3-methyl-            | 168 C13H12       | 000643-93-6 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

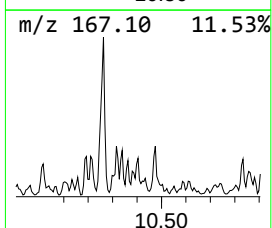
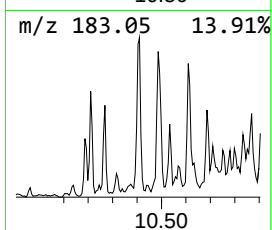
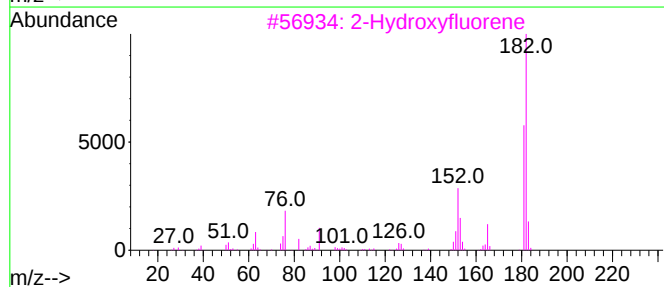
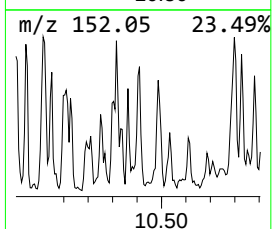
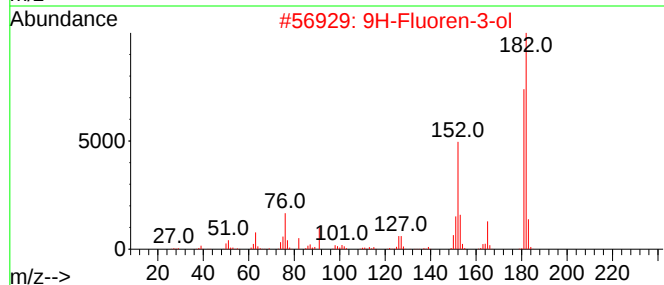
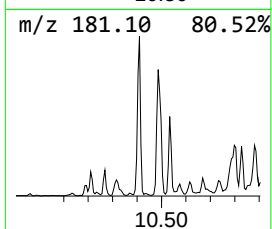
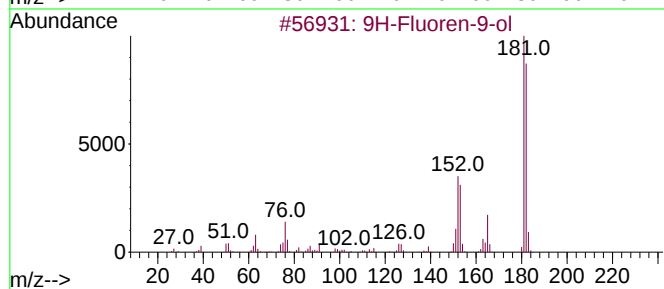
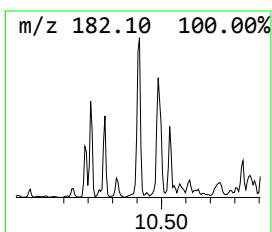
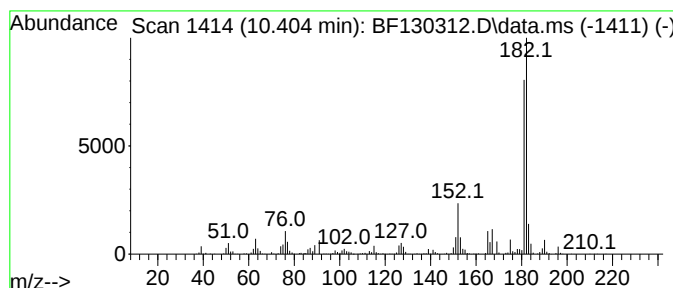
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 9H-Fluoren-9-ol Concentration Rank 18

| R.T.      | EstConc                 | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------|---------|------------------|-------------|------|
| 10.404    | 12.50 ng                | 1304780 | Acenaphthene-d10 | 9.704       |      |
| Hit# of 5 | Tentative ID            | MW      | MolForm          | CAS#        | Qual |
| 1         | 9H-Fluoren-9-ol         | 182     | C13H10O          | 001689-64-1 | 90   |
| 2         | 9H-Fluoren-3-ol         | 182     | C13H10O          | 006344-67-8 | 90   |
| 3         | 2-Hydroxyfluorene       | 182     | C13H10O          | 002443-58-5 | 87   |
| 4         | 2-Fluorenamine          | 181     | C13H11N          | 000153-78-6 | 86   |
| 5         | Dibenzofuran, 4-methyl- | 182     | C13H10O          | 007320-53-8 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

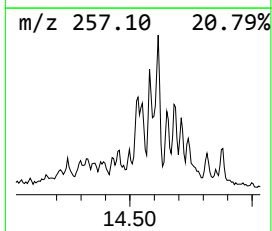
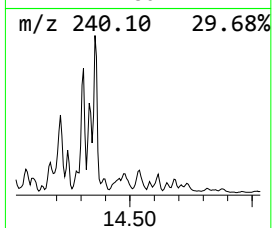
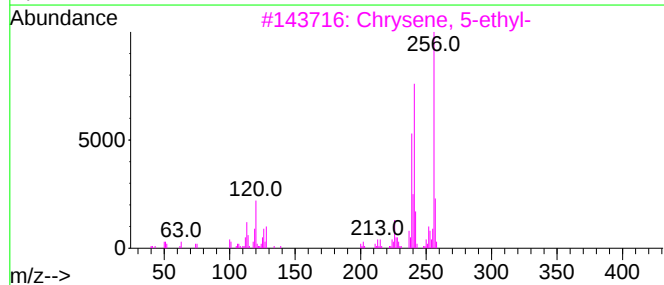
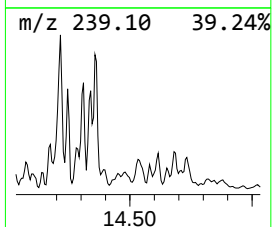
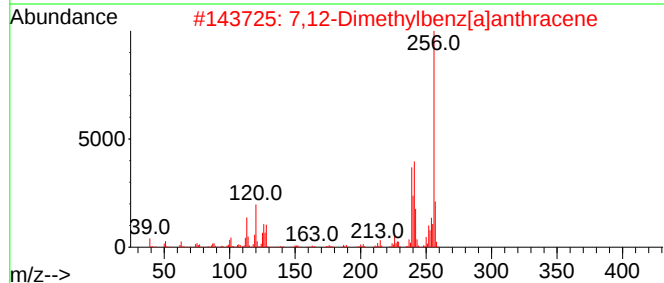
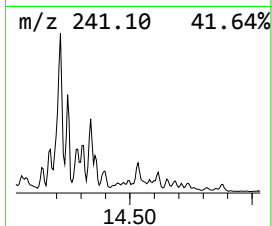
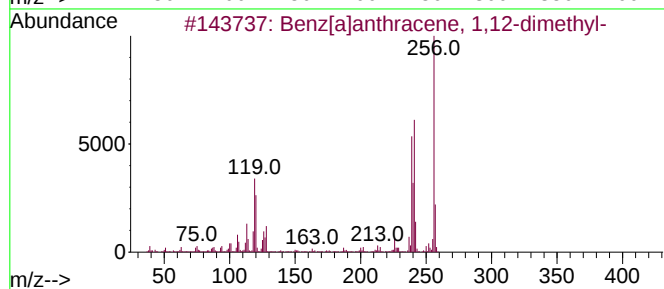
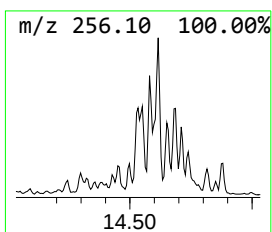
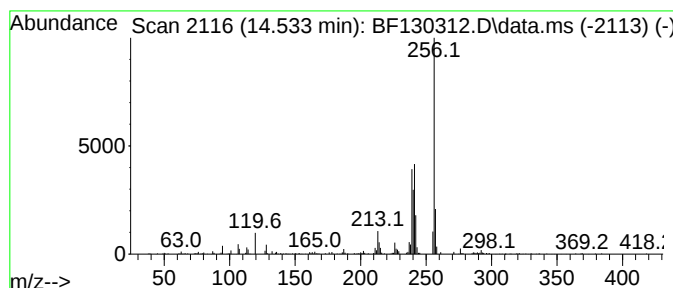
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 Benz[a]anthracene, 1,12-dim... Concentration Rank 10

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 14.533    | 21.99 ng                            | 850532 | Perylene-d12     | 15.221      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Benz[a]anthracene, 1,12-dimethyl-   | 256    | C20H16           | 000313-74-6 | 91   |
| 2         | 7,12-Dimethylbenz[a]anthracene      | 256    | C20H16           | 000057-97-6 | 91   |
| 3         | Chrysene, 5-ethyl-                  | 256    | C20H16           | 054986-62-8 | 91   |
| 4         | 5,12-Dimethylbenz[a]anthracene      | 256    | C20H16           | 021297-22-3 | 90   |
| 5         | Benzo[c]phenanthrene, 1,12-dimet... | 256    | C20H16           | 004076-43-1 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

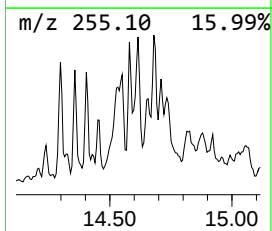
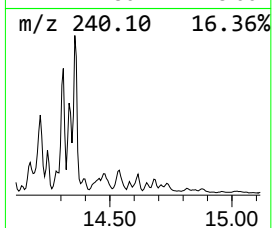
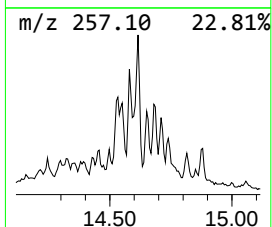
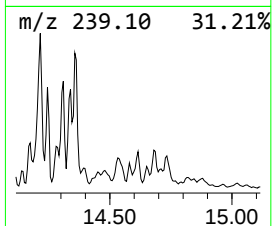
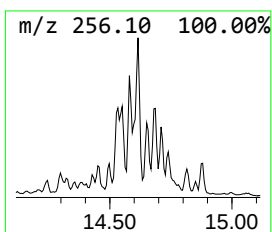
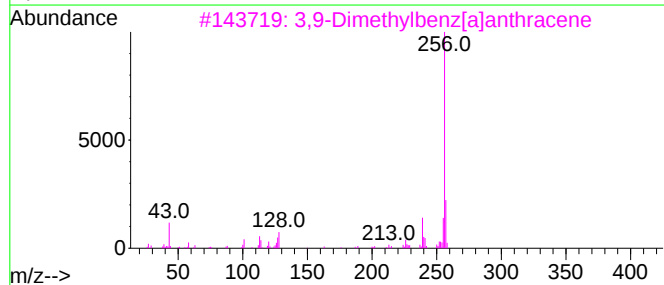
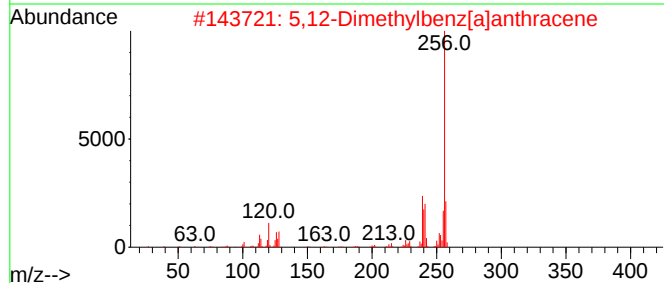
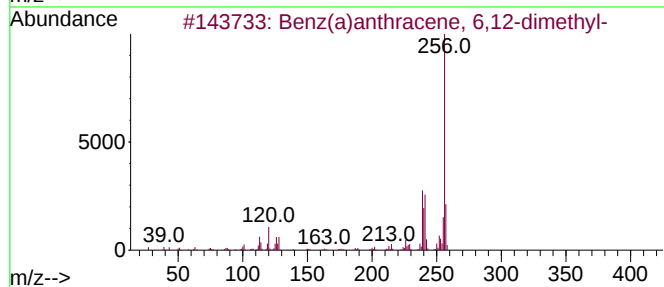
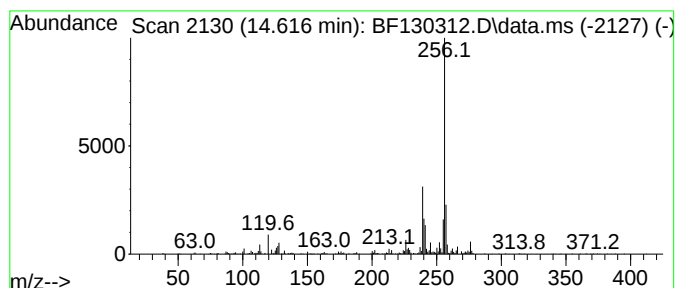
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 17 Benz(a)anthracene, 6,12-dim... Concentration Rank 15

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 14.616    | 17.43 ng                            | 674431 | Perylene-d12     | 15.221      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Benz(a)anthracene, 6,12-dimethyl-   | 256    | C20H16           | 000568-81-0 | 91   |
| 2         | 5,12-Dimethylbenz[a]anthracene      | 256    | C20H16           | 021297-22-3 | 91   |
| 3         | 3,9-Dimethylbenz[a]anthracene       | 256    | C20H16           | 000316-51-8 | 90   |
| 4         | Benz(a)anthracene, 8,12-dimethyl-   | 256    | C20H16           | 020627-31-0 | 87   |
| 5         | Benzo[c]phenanthrene, 5,8-dimethyl- | 256    | C20H16           | 054986-63-9 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

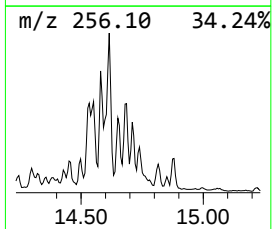
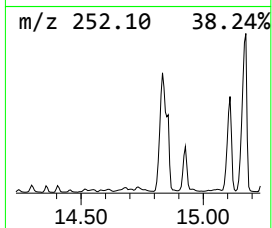
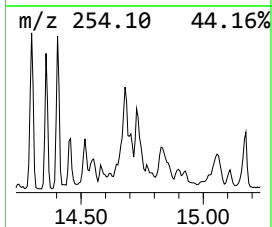
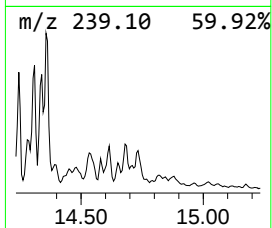
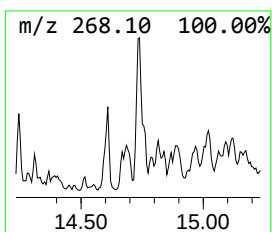
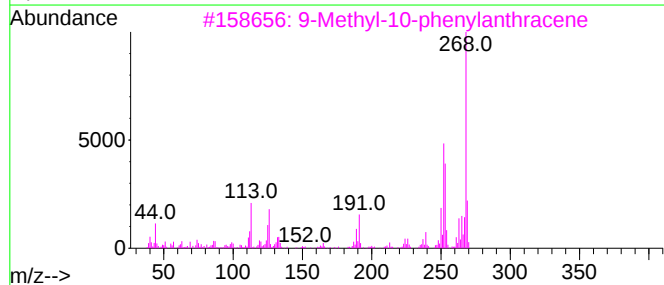
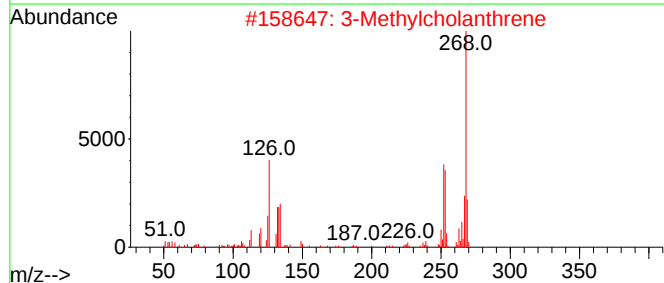
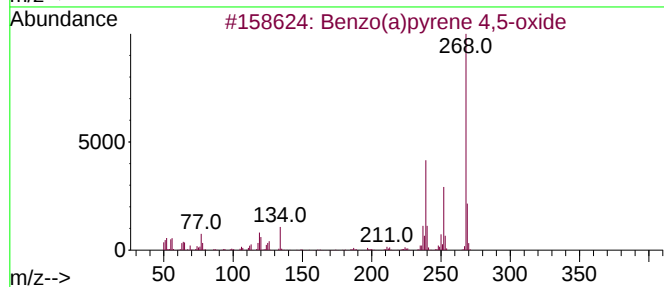
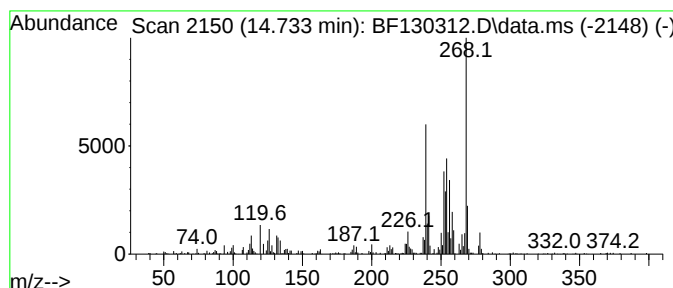
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 18 Benzo(a)pyrene 4,5-oxide Concentration Rank 9

| R.T.   | EstConc  | Area                                | Relative to ISTD | R.T.    |             |      |
|--------|----------|-------------------------------------|------------------|---------|-------------|------|
| 14.733 | 22.09 ng | 854441                              | Perylene-d12     | 15.221  |             |      |
| Hit#   | of 5     | Tentative ID                        | MW               | MolForm | CAS#        | Qual |
| 1      |          | Benzo(a)pyrene 4,5-oxide            | 268              | C20H12O | 037574-47-3 | 78   |
| 2      |          | 3-Methylcholanthrene                | 268              | C21H16  | 000056-49-5 | 55   |
| 3      |          | 9-Methyl-10-phenylanthracene        | 268              | C21H16  | 013425-08-6 | 43   |
| 4      |          | 9,10-Dihydro-8-methylbenzo[a]pyrene | 268              | C21H16  | 089524-72-1 | 41   |
| 5      |          | Dinaphtho[2,1-b:1',2'-d]furan       | 268              | C20H12O | 000194-63-8 | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

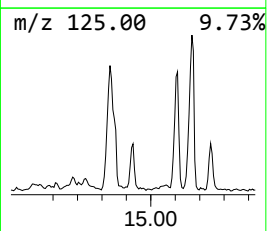
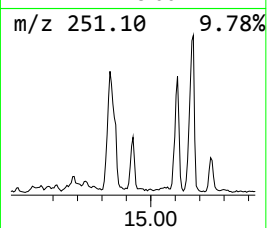
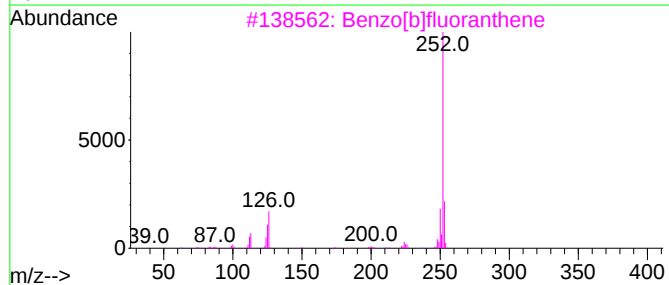
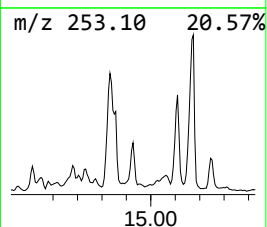
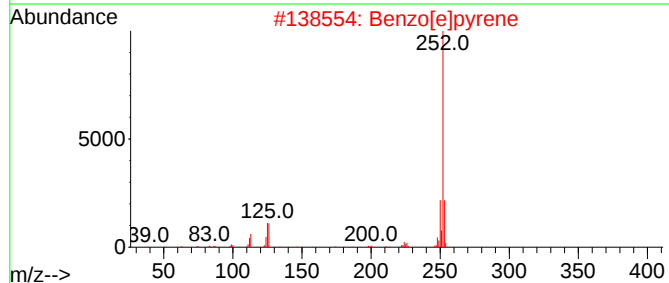
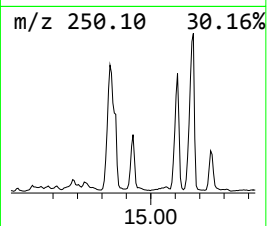
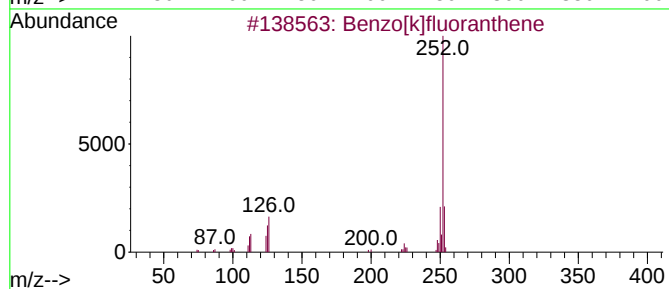
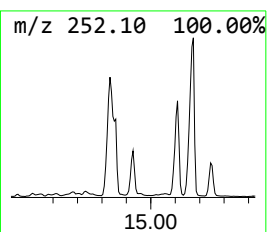
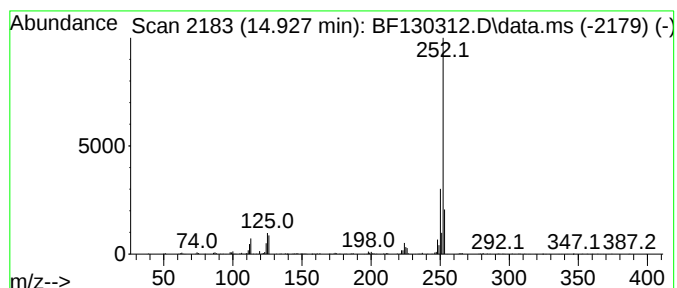
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 Benzo[e]pyrene Concentration Rank 16

| R.T.      | EstConc              | Area   | Relative to ISTD | R.T.        |      |
|-----------|----------------------|--------|------------------|-------------|------|
| 14.927    | 17.36 ng             | 671608 | Perylene-d12     | 15.221      |      |
| Hit# of 5 | Tentative ID         | MW     | MolForm          | CAS#        | Qual |
| 1         | Benzo[k]fluoranthene | 252    | C20H12           | 000207-08-9 | 96   |
| 2         | Benzo[e]pyrene       | 252    | C20H12           | 000192-97-2 | 94   |
| 3         | Benzo[b]fluoranthene | 252    | C20H12           | 000205-99-2 | 93   |
| 4         | Benzo[a]pyrene       | 252    | C20H12           | 000050-32-8 | 91   |
| 5         | Benzo[j]fluoranthene | 252    | C20H12           | 000205-82-3 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
Data File : BF130312.D  
Acq On : 17 Sep 2022 02:41  
Operator : CG\JU  
Sample : N4665-02 10X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
COMP-2

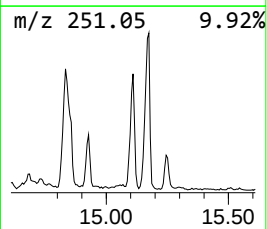
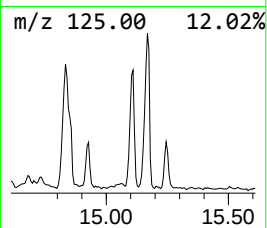
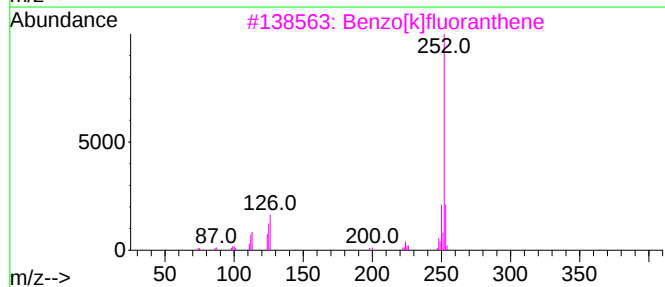
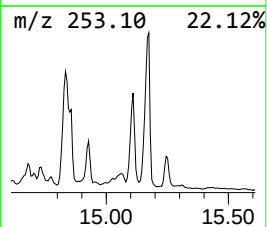
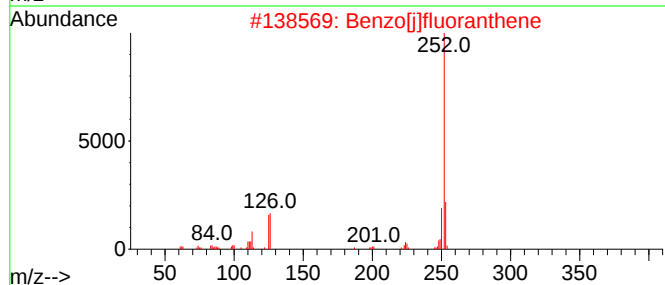
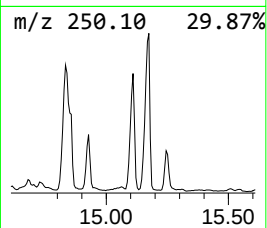
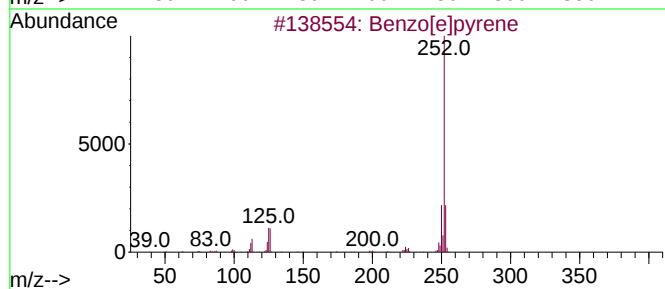
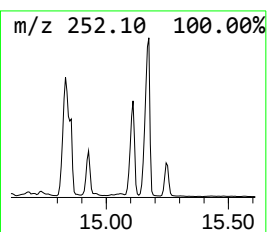
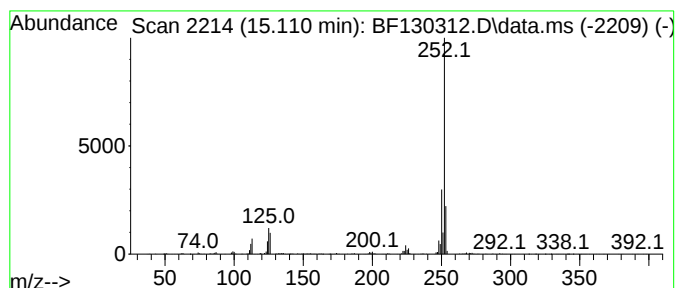
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 20 Benzo[j]fluoranthene Concentration Rank 5

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 15.110 | 40.03 ng | 1548420 | Perylene-d12     | 15.221 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------|-----|---------|-------------|------|
| 1    |      | Benzo[e]pyrene       | 252 | C20H12  | 000192-97-2 | 98   |
| 2    |      | Benzo[j]fluoranthene | 252 | C20H12  | 000205-82-3 | 94   |
| 3    |      | Benzo[k]fluoranthene | 252 | C20H12  | 000207-08-9 | 93   |
| 4    |      | Benzo[a]pyrene       | 252 | C20H12  | 000050-32-8 | 93   |
| 5    |      | Benzo[b]fluoranthene | 252 | C20H12  | 000205-99-2 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091622\  
 Data File : BF130312.D  
 Acq On : 17 Sep 2022 02:41  
 Operator : CG\JU  
 Sample : N4665-02 10X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 COMP-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Mesitylene         | 6.493  | 20.0    | ng    | 1267430  | 1                     | 6.669  | 1267430 | 20.0 |
| Indane             | 6.851  | 43.0    | ng    | 2725510  | 1                     | 6.669  | 1267430 | 20.0 |
| 1-Propyne, 3-ph... | 6.928  | 23.8    | ng    | 1510200  | 1                     | 6.669  | 1267430 | 20.0 |
| Naphthalene, 2-... | 9.222  | 22.2    | ng    | 2318260  | 3                     | 9.704  | 2088420 | 20.0 |
| Naphthalene, 1,... | 9.304  | 61.0    | ng    | 6366190  | 3                     | 9.704  | 2088420 | 20.0 |
| Naphthalene, 2,... | 9.375  | 67.8    | ng    | 7074470  | 3                     | 9.704  | 2088420 | 20.0 |
| Naphthalene, 1,... | 9.404  | 40.2    | ng    | 4197930  | 3                     | 9.704  | 2088420 | 20.0 |
| Naphthalene, 1,... | 9.486  | 36.7    | ng    | 3835870  | 3                     | 9.704  | 2088420 | 20.0 |
| Naphthalene, 2-... | 9.810  | 12.4    | ng    | 1297710  | 3                     | 9.704  | 2088420 | 20.0 |
| Naphthalene, 1,... | 10.022 | 20.0    | ng    | 2084830  | 3                     | 9.704  | 2088420 | 20.0 |
| Naphthalene, 1,... | 10.051 | 12.9    | ng    | 1349560  | 3                     | 9.704  | 2088420 | 20.0 |
| Naphthalene, 2,... | 10.110 | 18.5    | ng    | 1928740  | 3                     | 9.704  | 2088420 | 20.0 |
| 3,3'-Dimethylbi... | 10.210 | 18.7    | ng    | 1948030  | 3                     | 9.704  | 2088420 | 20.0 |
| 1-Isopropenylna... | 10.316 | 12.1    | ng    | 1267630  | 3                     | 9.704  | 2088420 | 20.0 |
| 9H-Fluoren-9-ol    | 10.404 | 12.5    | ng    | 1304780  | 3                     | 9.704  | 2088420 | 20.0 |
| Benz[a]anthrace... | 14.533 | 22.0    | ng    | 850532   | 6                     | 15.221 | 773682  | 20.0 |
| Benz(a)anthrace... | 14.616 | 17.4    | ng    | 674431   | 6                     | 15.221 | 773682  | 20.0 |
| Benzo(a)pyrene ... | 14.733 | 22.1    | ng    | 854441   | 6                     | 15.221 | 773682  | 20.0 |
| Benzo[e]pyrene     | 14.927 | 17.4    | ng    | 671608   | 6                     | 15.221 | 773682  | 20.0 |
| Benzo[j]fluoran... | 15.110 | 40.0    | ng    | 1548420  | 6                     | 15.221 | 773682  | 20.0 |