LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

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-BF073019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 5.457 | 570 | 573 | 594 | rBV | 1129171 | 1264500 | 25.73% | 1.581% |
| 2 | 6.469 | 742 | 745 | 757 | rBV | 1321296 | 1390133 | 28.29% | 1.738% |
| 3 | 6.610 | 765 | 769 | 776 | rBV | 1558438 | 1453766 | 29.59% | 1.817% |
| 4 | 6.669 | 776 | 779 | 784 | rVV | 224814 | 260879 | 5.31% | 0.326% |
| 5 | 6.834 | 804 | 807 | 815 | rBV | 978207 | 907676 | 18.47% | 1.135% |
| 6 | 6.992 | 830 | 834 | 849 | rVB | 1378678 | 1198071 | 24.38% | 1.498% |
| 7 | 7.398 | 900 | 903 | 909 | rBV | 870966 | 847370 | 17.25% | 1.059% |
| 8 | 7.692 | 951 | 953 | 962 | rVB | 238299 | 239753 | 4.88% | 0.300% |
| 9 | 8.116 | 1019 | 1025 | 1028 | rBV | 1393365 | 1227001 | 24.97% | 1.534% |
| 10 | 8.334 | 1057 | 1062 | 1068 | rBV | 226283 | 228559 | 4.65% | 0.286% |
| 11 | 8.392 | 1068 | 1072 | 1079 | rVB | 228548 | 248136 | 5.05% | 0.310% |
| 12 | 9.186 | 1202 | 1207 | 1213 | rBV | 2100509 | 1926544 | 39.21% | 2.408% |
| 13 | 9.528 | 1262 | 1265 | 1268 | rVV | 228648 | 243834 | 4.96% | 0.305% |
| 14 | 9.581 | 1271 | 1274 | 1281 | rVV4 | 564277 | 841422 | 17.12% | 1.052% |
| 15 | 9.733 | 1296 | 1300 | 1303 | rBV | 613553 | 603432 | 12.28% | 0.754% |
| 16 | 9.869 | 1317 | 1323 | 1327 | rBV | 1576338 | 1546808 | 31.48% | 1.934% |
| 17 | 10.075 | 1354 | 1358 | 1361 | rVV2 | 405229 | 448873 | 9.14% | 0.561% |
| 18 | 10.116 | 1361 | 1365 | 1368 | rVB | 521902 | 471312 | 9.59% | 0.589% |
| 19 | 10.186 | 1373 | 1377 | 1380 | rBV2 | 555139 | 667442 | 13.58% | 0.834% |
| 20 | 10.216 | 1380 | 1382 | 1384 | rVB | 318690 | 223071 | 4.54% | 0.279% |
| 21 | 10.281 | 1388 | 1393 | 1394 | rBV2 | 390540 | 450596 | 9.17% | 0.563% |
| 22 | 10.298 | 1394 | 1396 | 1398 | rVB2 | 307389 | 235826 | 4.80% | 0.295% |
| 23 | 10.469 | 1422 | 1425 | 1429 | rBV3 | 142745 | 241333 | 4.91% | 0.302% |
| 24 | 10.528 | 1433 | 1435 | 1439 | rVB3 | 260957 | 251579 | 5.12% | 0.314% |
| 25 | 10.663 | 1455 | 1458 | 1461 | rVV | 1038162 | 1063026 | 21.63% | 1.329% |
| 26 | 10.786 | 1474 | 1479 | 1484 | rVB2 | 792499 | 919176 | 18.71% | 1.149% |
| 27 | 10.857 | 1488 | 1491 | 1494 | rBV | 353866 | 343511 | 6.99% | 0.429% |
| 28 | 10.969 | 1505 | 1510 | 1512 | rVV3 | 321487 | 484029 | 9.85% | 0.605% |
| 29 | 10.998 | 1512 | 1515 | 1517 | rVV | 358390 | 401257 | 8.17% | 0.502% |
| 30 | 11.022 | 1517 | 1519 | 1522 | rVV3 | 194426 | 294643 | 6.00% | 0.368% |
| 31 | 11.098 | 1526 | 1532 | 1534 | rBV3 | 253140 | 451547 | 9.19% | 0.564% |
| 32 | 11.116 | 1534 | 1535 | 1540 | rVV3 | 287464 | 335849 | 6.83% | 0.420% |
| 33 | 11.169 | 1540 | 1544 | 1546 | rVV4 | 271824 | 344274 | 7.01% | 0.430% |
| 34 | 11.216 | 1548 | 1552 | 1555 | rVV3 | 247735 | 397690 | 8.09% | 0.497% |

LSC Area Percent Report

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Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

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-BF073019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 11.251 | 1555 | 1558 | 1564 | rVB3 | 392751 | 494453 | 10.06% | 0.618% |
| 36 | 11.357 | 1573 | 1576 | 1578 | rBV | 1335911 | 1216041 | 24.75% | 1.520% |
| 37 | 11.380 | 1578 | 1580 | 1587 | rVB | 1612682 | 1469520 | 29.91% | 1.837% |
| 38 | 11.433 | 1587 | 1589 | 1592 | rBV | 582609 | 446142 | 9.08% | 0.558% |
| 39 | 11.469 | 1592 | 1595 | 1598 | rBV3 | 285729 | 376315 | 7.66% | 0.470% |
| 40 | 11.598 | 1613 | 1617 | 1622 | rBV6 | 204526 | 485375 | 9.88% | 0.607% |
| 41 | 11.645 | 1623 | 1625 | 1628 | rVV3 | 374918 | 387255 | 7.88% | 0.484% |
| 42 | 11.692 | 1628 | 1633 | 1638 | rVV2 | 521529 | 685063 | 13.94% | 0.856% |
| 43 | 11.775 | 1644 | 1647 | 1650 | rVB | 310300 | 339496 | 6.91% | 0.424% |
| 44 | 11.863 | 1658 | 1662 | 1664 | rBV | 1434574 | 1317761 | 26.82% | 1.647% |
| 45 | 11.892 | 1664 | 1667 | 1671 | rVB | 1903634 | 1647459 | 33.53% | 2.059% |
| 46 | 11.939 | 1672 | 1675 | 1677 | rBV | 496276 | 505272 | 10.28% | 0.632% |
| 47 | 11.975 | 1677 | 1681 | 1683 | rVV | 1598543 | 1768739 | 36.00% | 2.211% |
| 48 | 11.998 | 1683 | 1685 | 1689 | rVB | 1131969 | 891754 | 18.15% | 1.115% |
| 49 | 12.092 | 1699 | 1701 | 1704 | rBV | 360702 | 307771 | 6.26% | 0.385% |
| 50 | 12.151 | 1708 | 1711 | 1714 | rBV2 | 648863 | 724272 | 14.74% | 0.905% |
| 51 | 12.186 | 1716 | 1717 | 1722 | rVB2 | 471167 | 461846 | 9.40% | 0.577% |
| 52 | 12.304 | 1732 | 1737 | 1740 | rBV2 | 720198 | 938480 | 19.10% | 1.173% |
| 53 | 12.339 | 1740 | 1743 | 1745 | rVV | 942804 | 807586 | 16.44% | 1.009% |
| 54 | 12.363 | 1745 | 1747 | 1750 | rVB | 710055 | 542808 | 11.05% | 0.679% |
| 55 | 12.416 | 1752 | 1756 | 1758 | rBV | 1914396 | 1996065 | 40.62% | 2.495% |
| 56 | 12.451 | 1758 | 1762 | 1764 | rVV3 | 920177 | 1301085 | 26.48% | 1.626% |
| 57 | 12.504 | 1768 | 1771 | 1775 | rBV3 | 693727 | 991704 | 20.18% | 1.240% |
| 58 | 12.580 | 1780 | 1784 | 1787 | rBV | 2354764 | 2287648 | 46.56% | 2.860% |
| 59 | 12.616 | 1787 | 1790 | 1792 | rBV2 | 369495 | 318242 | 6.48% | 0.398% |
| 60 | 12.639 | 1792 | 1794 | 1797 | rVB | 400817 | 317936 | 6.47% | 0.397% |
| 61 | 12.733 | 1806 | 1810 | 1812 | rBV4 | 240183 | 298459 | 6.07% | 0.373% |
| 62 | 12.757 | 1812 | 1814 | 1817 | rVB2 | 340452 | 262487 | 5.34% | 0.328% |
| 63 | 12.816 | 1818 | 1824 | 1828 | rBV | 3629369 | 4913681 | 100.00% | 6.142% |
| 64 | 12.863 | 1828 | 1832 | 1835 | rVB2 | 974159 | 1019367 | 20.75% | 1.274% |
| 65 | 12.916 | 1835 | 1841 | 1842 | rBV2 | 369812 | 598548 | 12.18% | 0.748% |
| 66 | 12.939 | 1842 | 1845 | 1848 | rBV | 1659314 | 1248119 | 25.40% | 1.560% |
| 67 | 12.963 | 1848 | 1849 | 1855 | rVB2 | 438325 | 391206 | 7.96% | 0.489% |
| 68 | 13.022 | 1855 | 1859 | 1866 | rBV4 | 794082 | 1415403 | 28.81% | 1.769% |
| 69 | 13.080 | 1866 | 1869 | 1871 | rBV | 320443 | 312458 | 6.36% | 0.391% |
| 70 | 13.116 | 1871 | 1875 | 1876 | rBV2 | 653064 | 785490 | 15.99% | 0.982% |
| 71 | 13.239 | 1893 | 1896 | 1900 | rVB | 1431474 | 1363947 | 27.76% | 1.705% |

LSC Area Percent Report

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Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

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-BF073019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 13.333 | 1908 | 1912 | 1914 | rVV | 1332208 | 1205193 | 24.53% | 1.507% |
| 73 | 13.363 | 1914 | 1917 | 1920 | rVV | 1289865 | 1145482 | 23.31% | 1.432% |
| 74 | 13.392 | 1920 | 1922 | 1926 | rVB4 | 315400 | 320029 | 6.51% | 0.400% |
| 75 | 13.445 | 1929 | 1931 | 1934 | rVB | 247017 | 230303 | 4.69% | 0.288% |
| 76 | 13.645 | 1962 | 1965 | 1969 | rVB | 653596 | 683231 | 13.90% | 0.854% |
| 77 | 13.704 | 1972 | 1975 | 1978 | rVB | 435866 | 351233 | 7.15% | 0.439% |
| 78 | 13.757 | 1978 | 1984 | 1986 | rBV3 | 641433 | 1172222 | 23.86% | 1.465% |
| 79 | 13.810 | 1990 | 1993 | 1996 | rBV2 | 323382 | 512845 | 10.44% | 0.641% |
| 80 | 13.857 | 1999 | 2001 | 2004 | rVB | 463086 | 400710 | 8.15% | 0.501% |
| 81 | 13.998 | 2021 | 2025 | 2029 | rBV2 | 1922308 | 2996378 | 60.98% | 3.746% |
| 82 | 14.033 | 2029 | 2031 | 2034 | rVV | 1699829 | 1345252 | 27.38% | 1.682% |
| 83 | 14.086 | 2038 | 2040 | 2043 | rBV | 347421 | 384470 | 7.82% | 0.481% |
| 84 | 14.321 | 2078 | 2080 | 2083 | rVB2 | 228776 | 205477 | 4.18% | 0.257% |
| 85 | 14.351 | 2083 | 2085 | 2087 | rBV | 322197 | 299050 | 6.09% | 0.374% |
| 86 | 14.392 | 2087 | 2092 | 2095 | rVV | 1103553 | 1180765 | 24.03% | 1.476% |
| 87 | 14.427 | 2095 | 2098 | 2100 | rVB | 476379 | 417577 | 8.50% | 0.522% |
| 88 | 14.463 | 2101 | 2104 | 2107 | rBV3 | 256281 | 321028 | 6.53% | 0.401% |
| 89 | 14.516 | 2111 | 2113 | 2115 | rBV | 442809 | 386117 | 7.86% | 0.483% |
| 90 | 14.580 | 2122 | 2124 | 2128 | rVB2 | 308399 | 320599 | 6.52% | 0.401% |
| 91 | 14.721 | 2145 | 2148 | 2150 | rBV | 230208 | 274311 | 5.58% | 0.343% |
| 92 | 14.886 | 2173 | 2176 | 2179 | rBV2 | 251325 | 251328 | 5.11% | 0.314% |
| 93 | 15.051 | 2200 | 2204 | 2210 | rBV2 | 856975 | 1406882 | 28.63% | 1.759% |
| 94 | 15.157 | 2219 | 2222 | 2226 | rVB | 192894 | 223505 | 4.55% | 0.279% |
| 95 | 15.351 | 2251 | 2255 | 2260 | rVB | 917587 | 1109702 | 22.58% | 1.387% |
| 96 | 15.415 | 2261 | 2266 | 2270 | rVV | 1194134 | 1564935 | 31.85% | 1.956% |
| 97 | 15.474 | 2272 | 2276 | 2279 | rVV | 681642 | 894031 | 18.19% | 1.118% |
| 98 | 15.504 | 2279 | 2281 | 2288 | rVB2 | 250311 | 406865 | 8.28% | 0.509% |
| 99 | 16.945 | 2521 | 2526 | 2534 | rVB | 421584 | 812176 | 16.53% | 1.015% |
| 100 | 17.392 | 2596 | 2602 | 2611 | rVB | 526870 | 1089444 | 22.17% | 1.362% |

Sum of corrected areas: 79999311

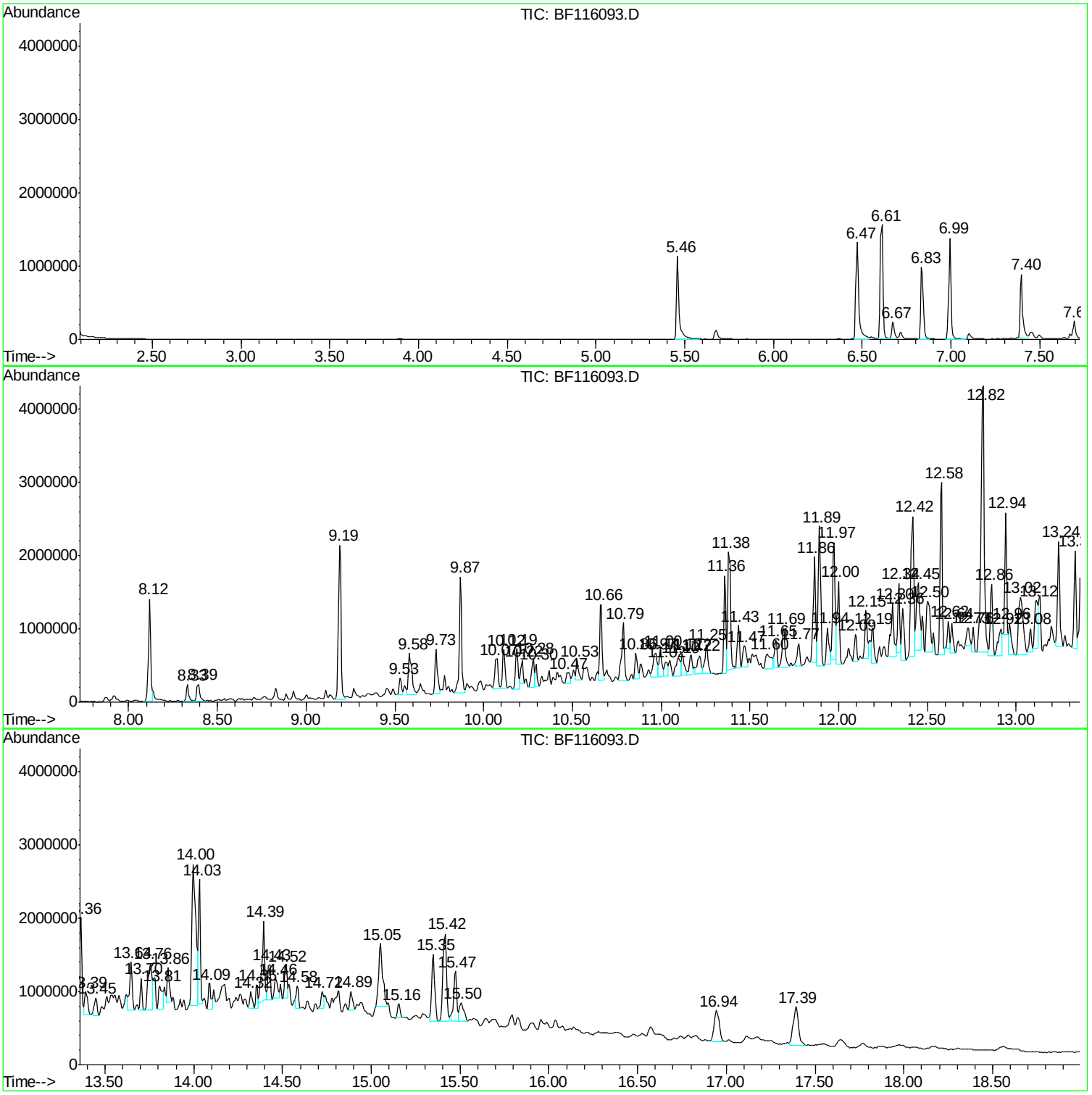
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

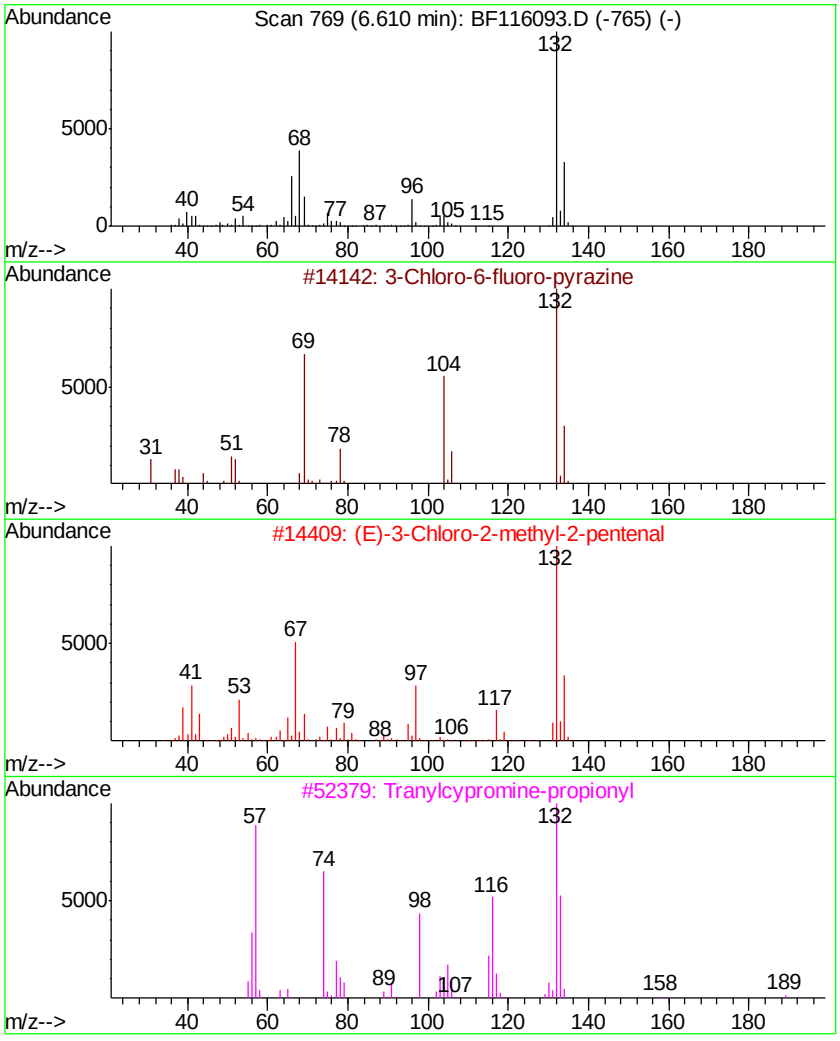
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

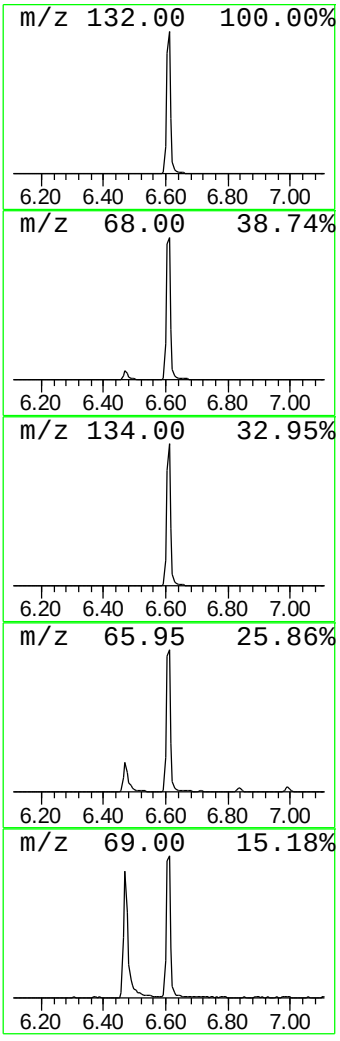
**********

Peak Number 1 unknown6.61 Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.61 | 32.03 ng | 1453770 | 1,4-Dichlorobenzene-d4 | 6.83 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Chloro-6-fluoro-pyrazine | 132 | C4H2ClFN2 | 1000146-10-7 | 16 |
| 2 | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO | 031357-76-3 | 16 |
| 3 | Tranylcypromine-propionyl | 189 | C12H15NO | 1000123-86-3 | 12 |
| 4 | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12 | 028749-81-7 | 9 |
| 5 | 4-Chloro-2-fluoro-pyrimidine | 132 | C4H2ClFN2 | 051422-00-5 | 9 |



Library Search Compound Report

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Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
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**Instrument :** BNA_F
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

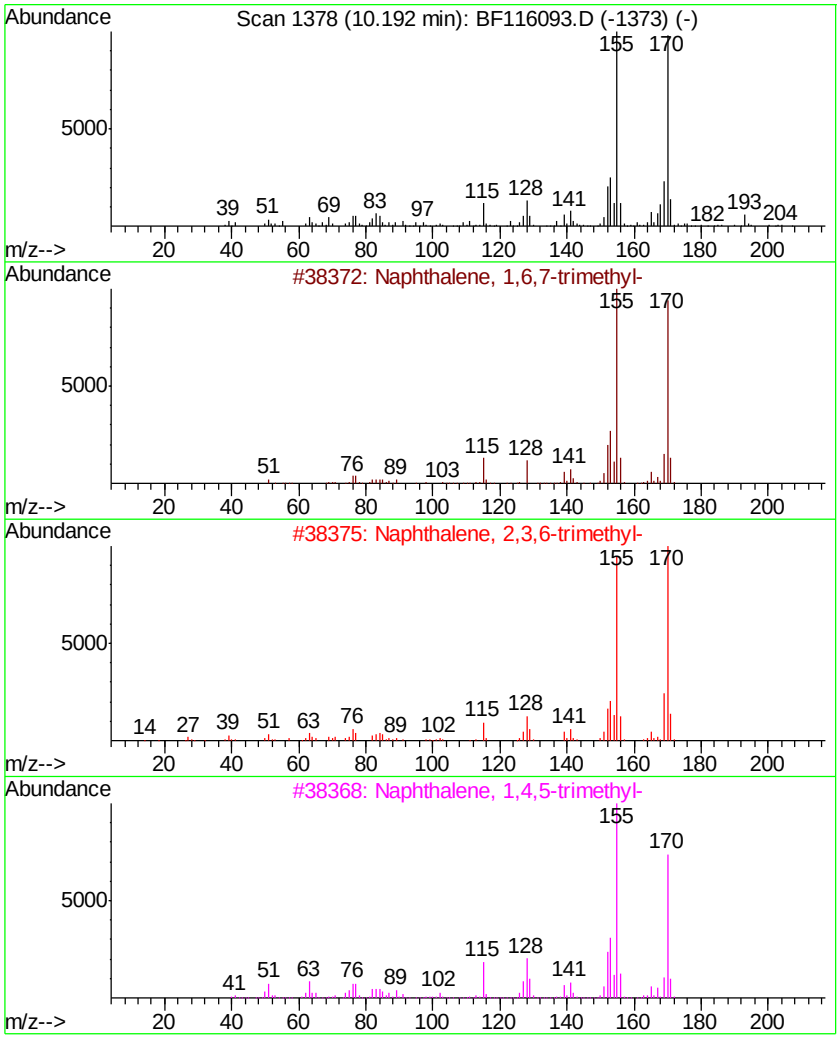
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

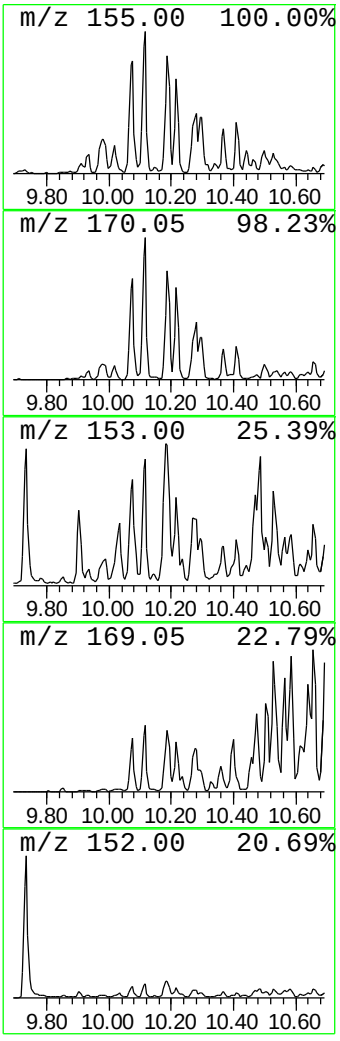
**********

Peak Number 3 Naphthalene, 1,6,7-trimethyl- Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.19 | 8.63 ng | 667442 | Acenaphthene-d10 | 9.87 |

| 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 | 96 |
| 5 | 4,6,8-Trimethylazulene | 170 | C13H14 | 000941-81-1 | 90 |



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Sample : K4178-01 2X
Misc :
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**Instrument :** BNA_F
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

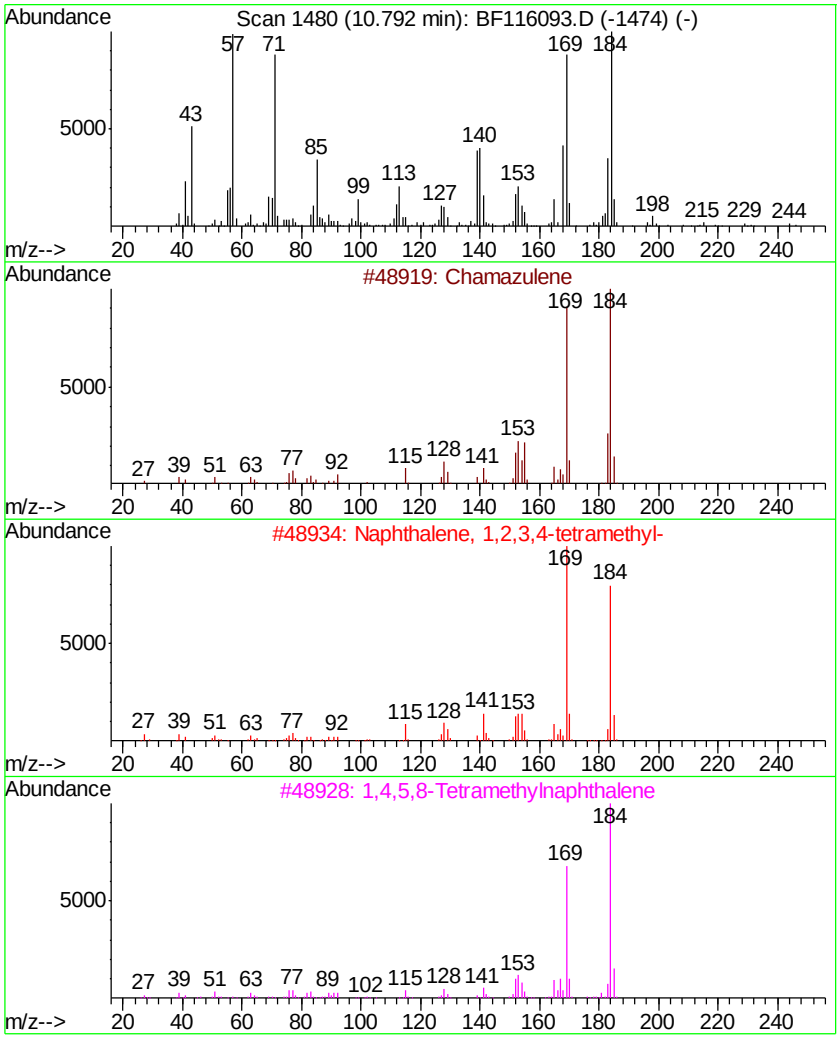
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

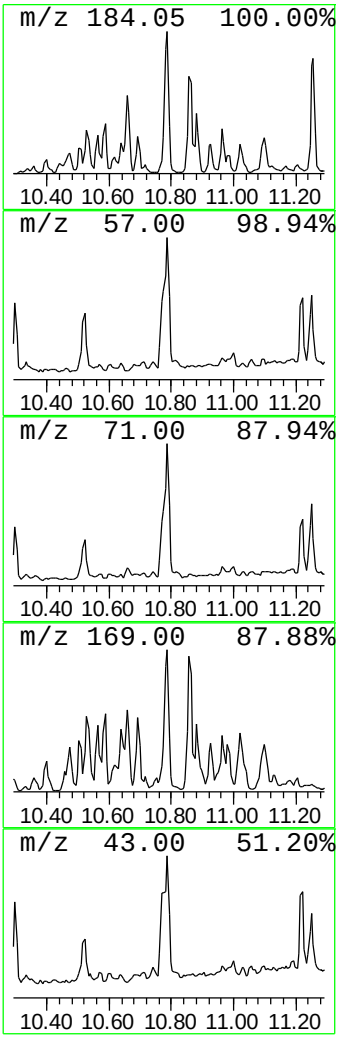
**********

Peak Number 4 Chamazulene Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.79 | 15.12 ng | 919176 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Chamazulene | 184 | C14H16 | 000529-05-5 | 93 |
| 2 | Naphthalene, 1,2,3,4-tetramethyl- | 184 | C14H16 | 003031-15-0 | 83 |
| 3 | 1,4,5,8-Tetramethylnaphthalene | 184 | C14H16 | 002717-39-7 | 64 |
| 4 | Cyclopentene, 1,2-dimethyl-4-met... | 184 | C14H16 | 1000152-22-4 | 64 |
| 5 | 3,4,4-Trimethyl-5-methylene-2-et... | 184 | C9H16N2S | 037120-36-8 | 52 |



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Sample : K4178-01 2X
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**Instrument :** BNA_F
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

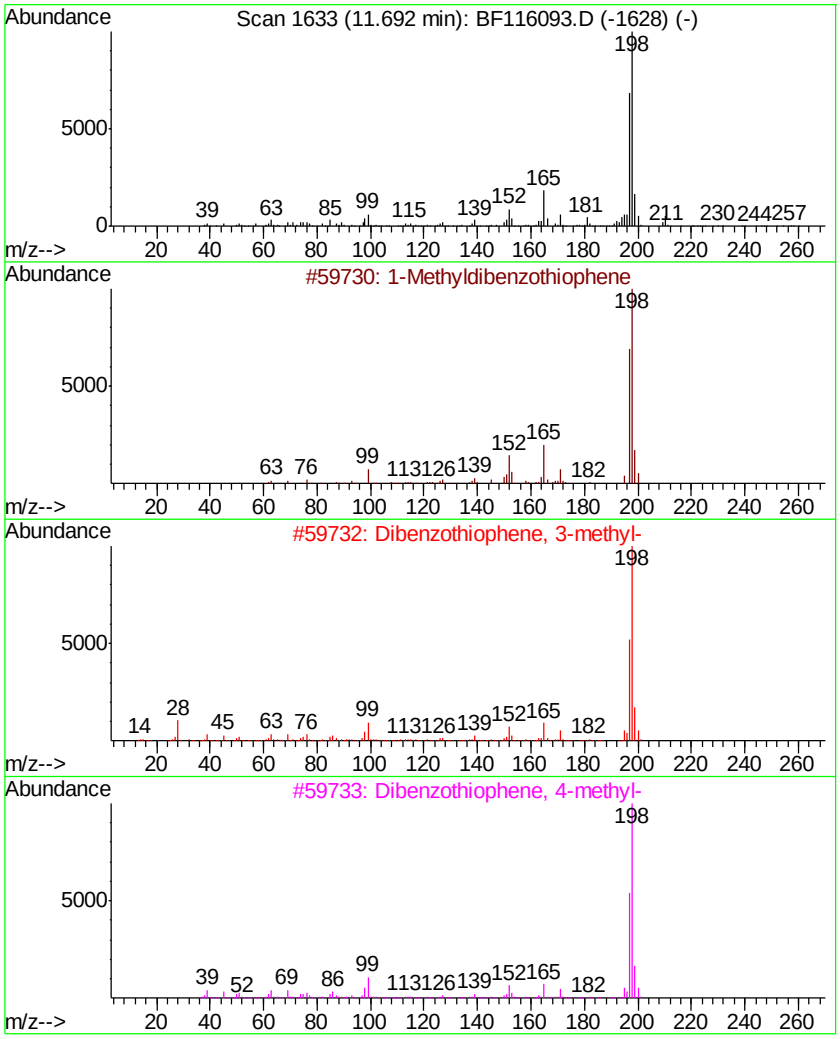
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

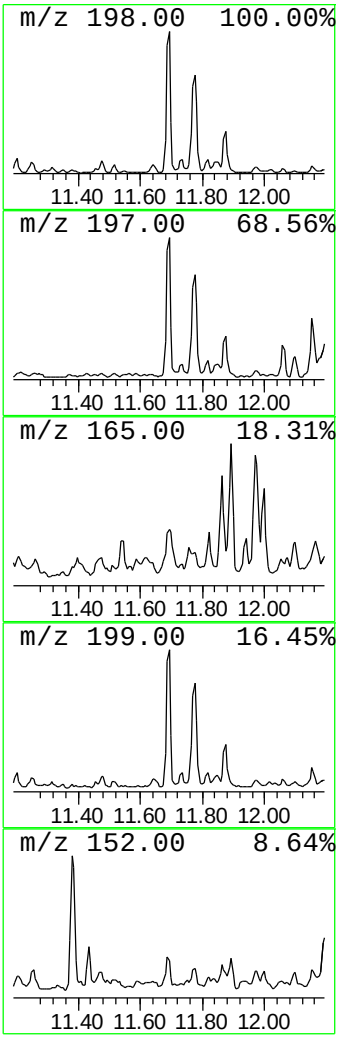
**********

Peak Number 5 1-Methyldibenzothiophene Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.69 | 11.27 ng | 685063 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Methyldibenzothiophene | 198 | C13H10S | 031317-07-4 | 93 |
| 2 | Dibenzothiophene, 3-methyl- | 198 | C13H10S | 016587-52-3 | 93 |
| 3 | Dibenzothiophene, 4-methyl- | 198 | C13H10S | 007372-88-5 | 93 |
| 4 | Benzenamine, 4,4'-methylenebis- | 198 | C13H14N2 | 000101-77-9 | 72 |
| 5 | Benzene, 1-fluoro-4-(2-phenyleth... | 198 | C14H11F | 017404-69-2 | 72 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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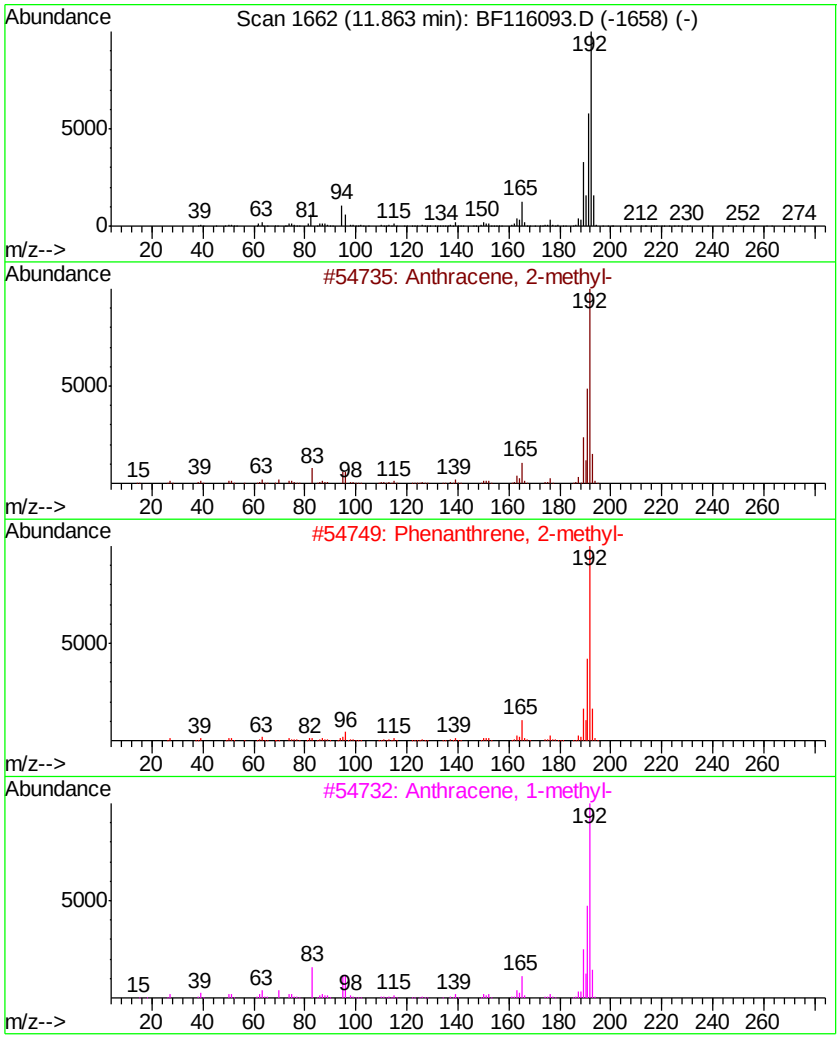
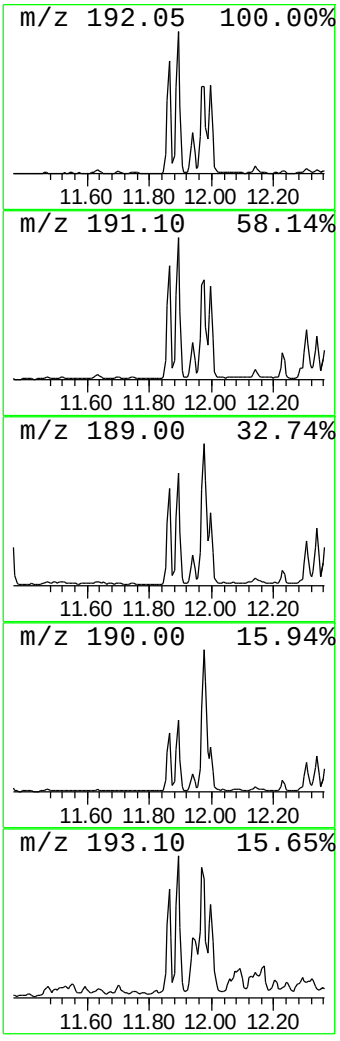
Peak Number 6 Anthracene, 2-methyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.86 | 21.67 ng | 1317760 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 95 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 92 |
| 3 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 91 |
| 4 | Anthracene, 9-methyl- | 192 | C15H12 | 000779-02-2 | 91 |
| 5 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 91 |



Library Search Compound Report

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Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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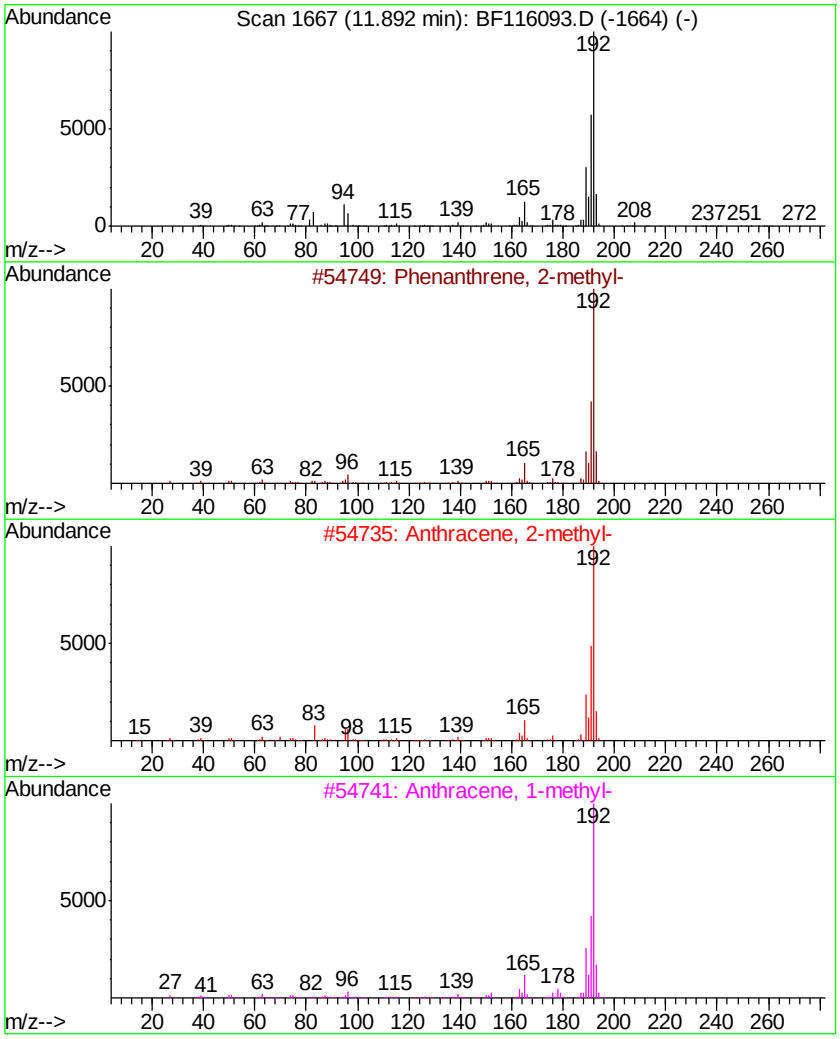
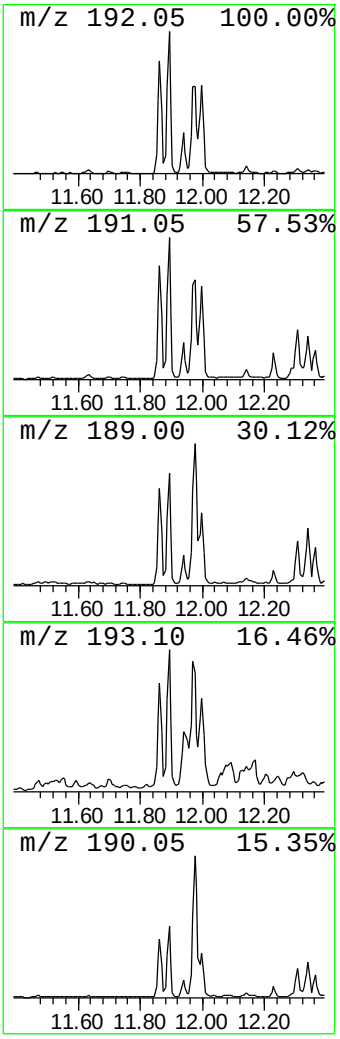
Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.89 | 27.10 ng | 1647460 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 96 |
| 2 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 96 |
| 3 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 95 |
| 4 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 93 |
| 5 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 93 |



Library Search Compound Report

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Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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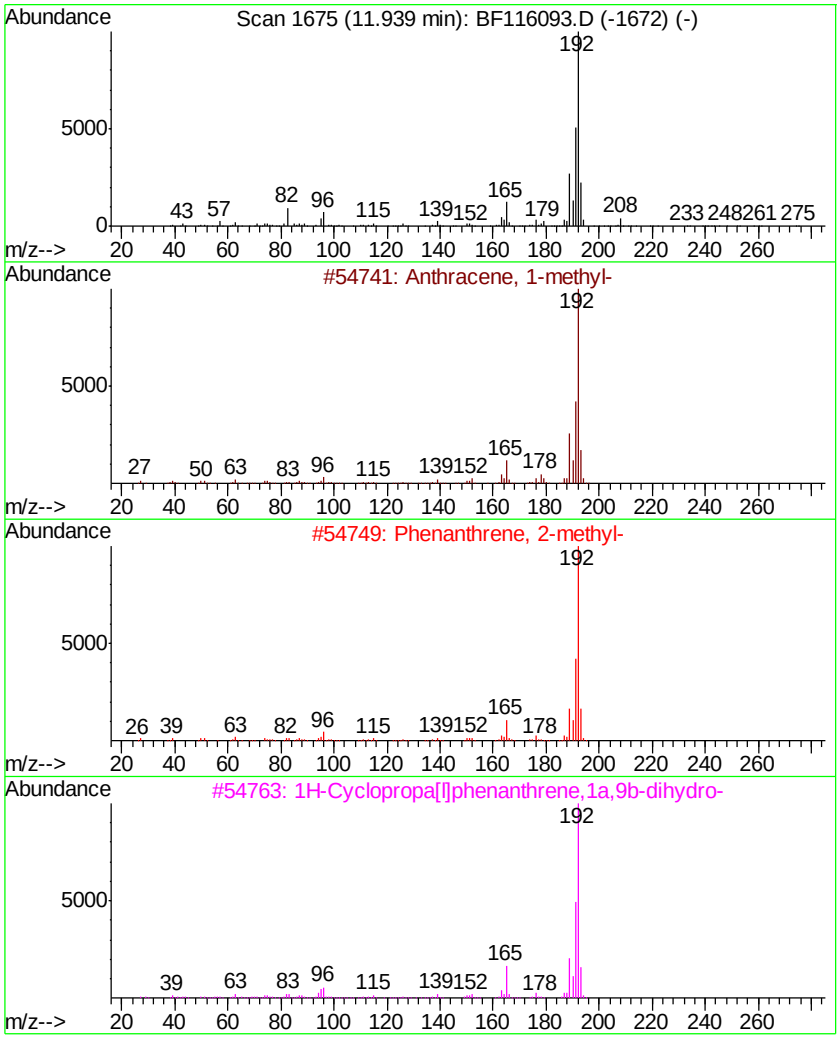
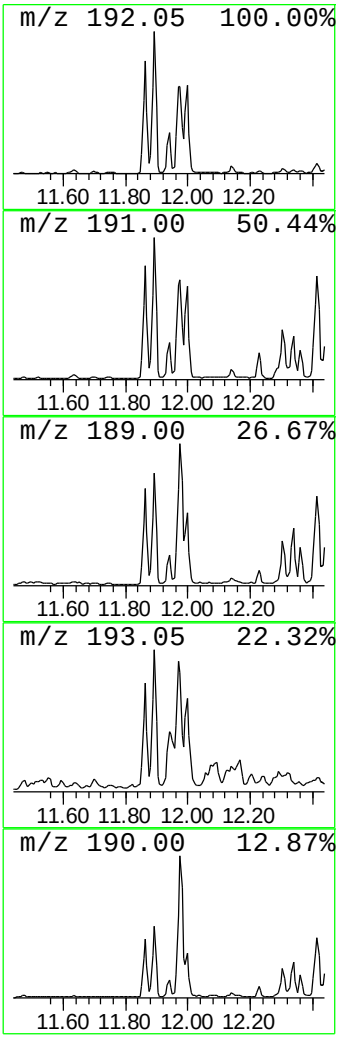
Peak Number 8 Anthracene, 1-methyl- Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.94 | 8.31 ng | 505272 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 96 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 96 |
| 3 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 95 |
| 4 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 93 |
| 5 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 90 |



# Library Search Compound Report

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Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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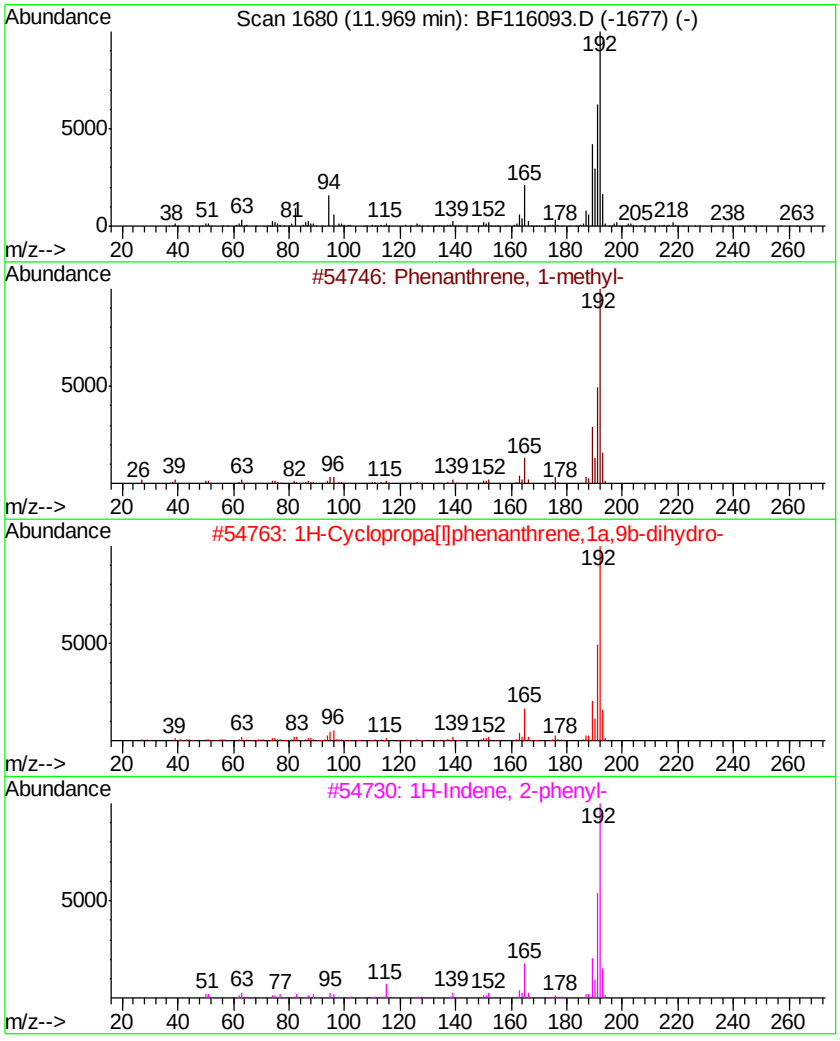
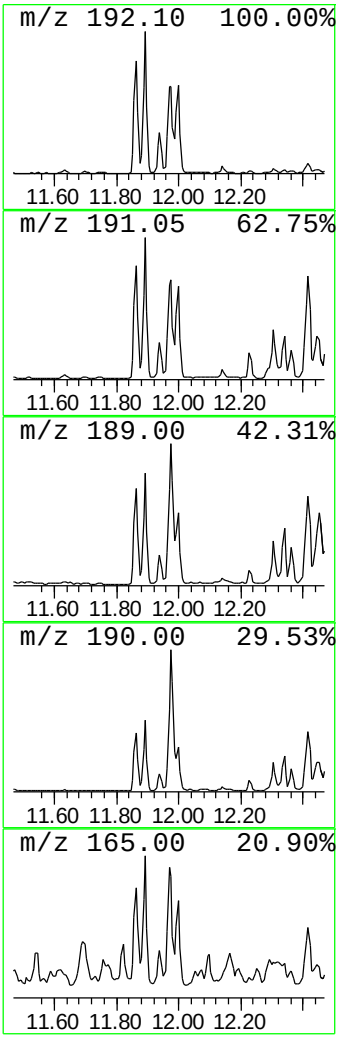
Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.97 | 29.09 ng | 1768740 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 60 |
| 2 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 60 |
| 3 | 1H-Indene, 2-phenyl- | 192 | C15H12 | 004505-48-0 | 55 |
| 4 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 55 |
| 5 | Phenanthrene, 4-methyl- | 192 | C15H12 | 000832-64-4 | 53 |



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Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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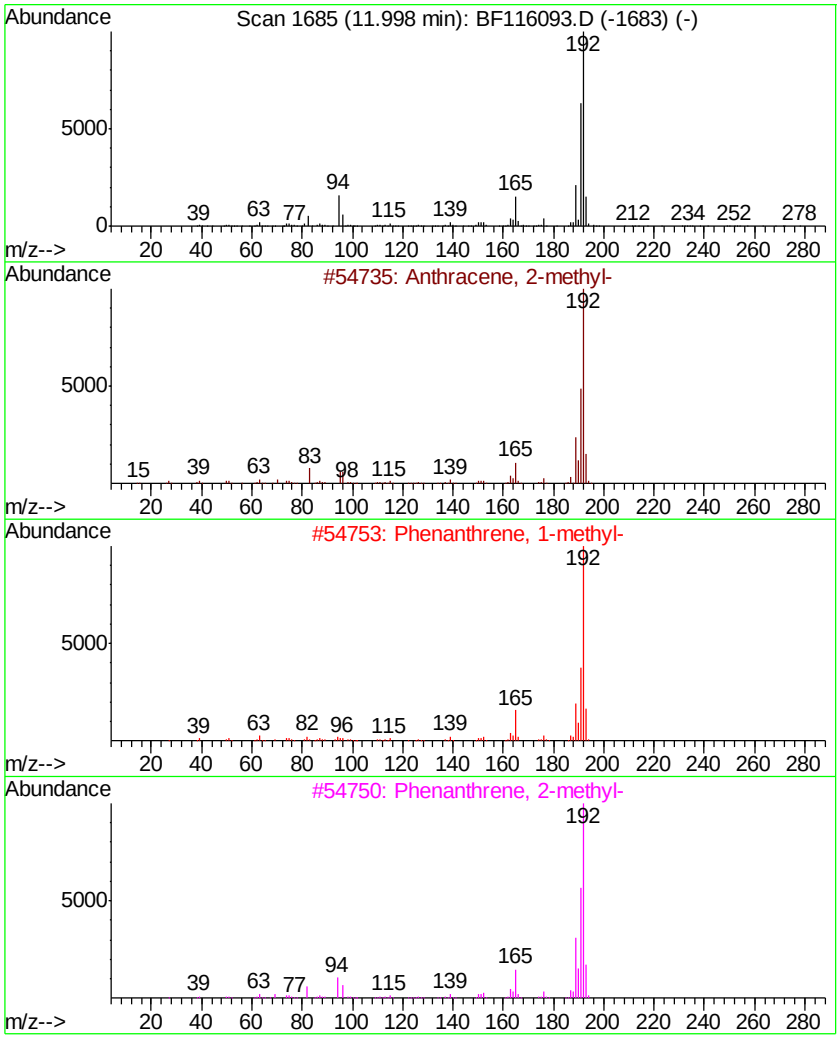
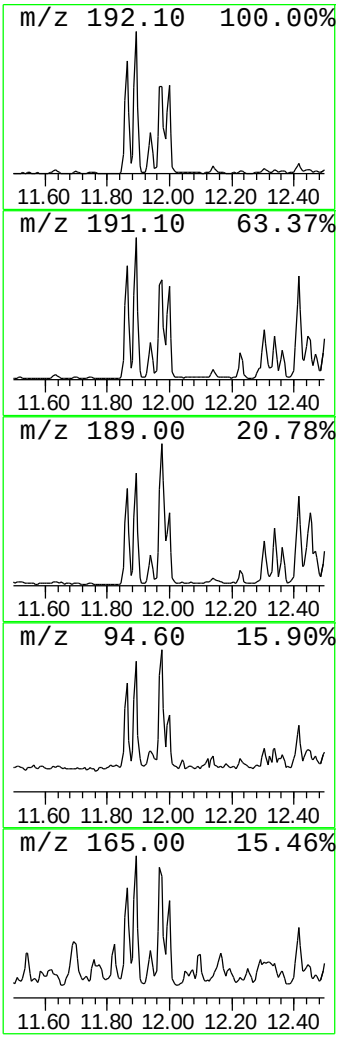
Peak Number 10 1H-Indene, 2-phenyl- Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.00 | 14.67 ng | 891754 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 95 |
| 2 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 92 |
| 3 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 91 |
| 4 | 1H-Indene, 2-phenyl- | 192 | C15H12 | 004505-48-0 | 90 |
| 5 | Propyne, 1,3-diphenyl- | 192 | C15H12 | 004980-70-5 | 81 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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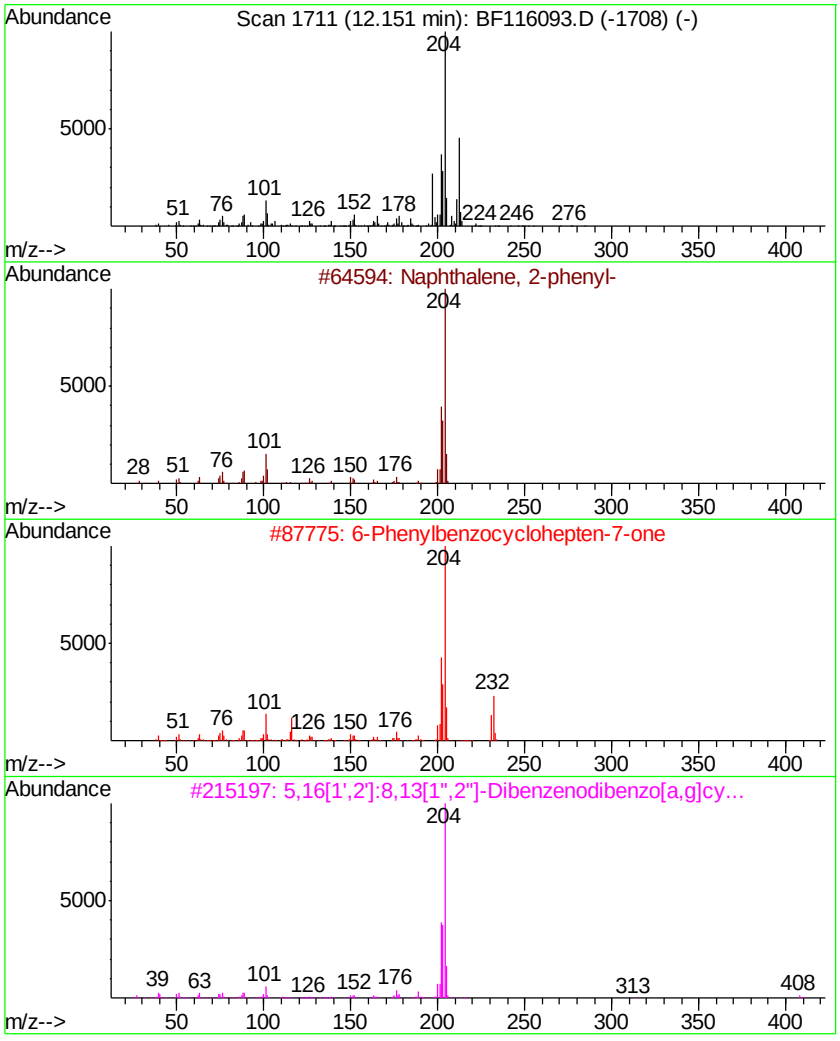
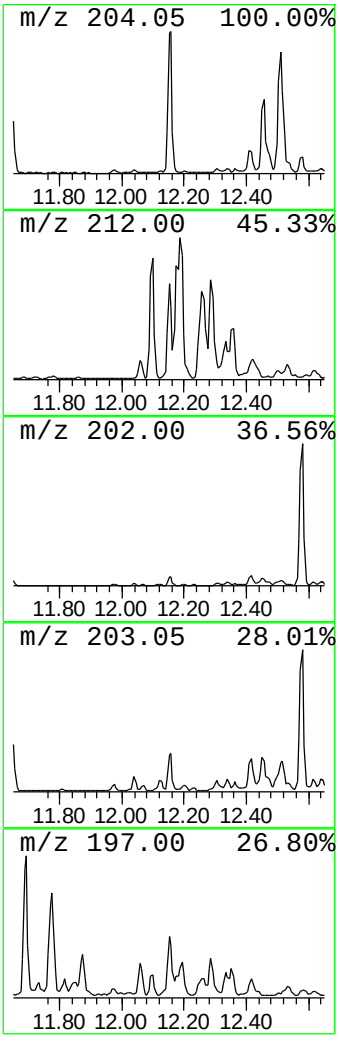
Peak Number 11 Naphthalene, 2-phenyl- Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.15 | 11.91 ng | 724272 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 64 |
| 2 | 6-Phenylbenzocyclohepten-7-one | 232 | C17H12O | 093327-56-1 | 49 |
| 3 | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24 | 005672-97-9 | 47 |
| 4 | 6-Cyanophenanthridine | 204 | C14H8N2 | 002176-28-5 | 41 |
| 5 | [1,1'-Biphenyl]-4,4'-dicarbonitrile | 204 | C14H8N2 | 001591-30-6 | 38 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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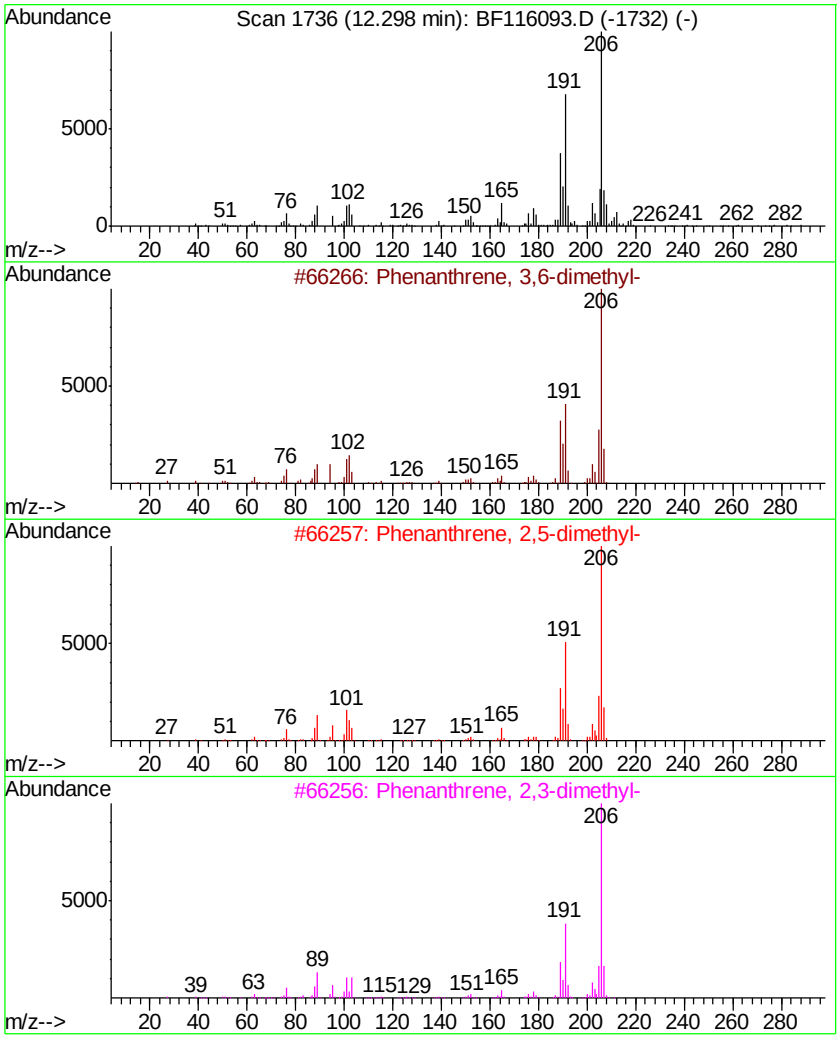
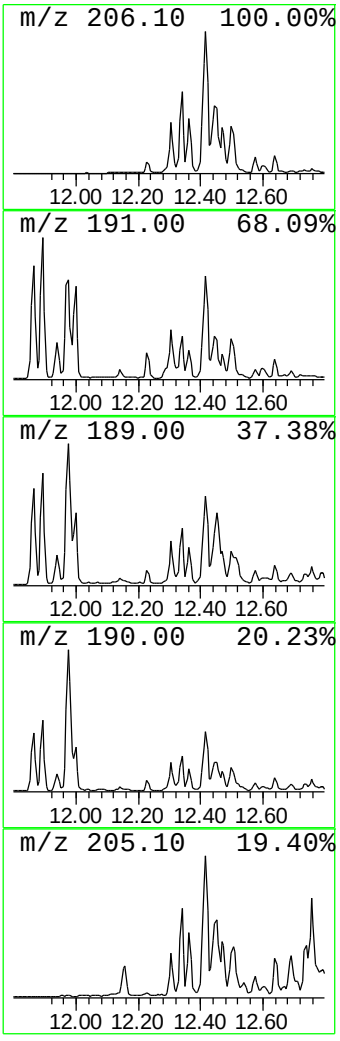
Peak Number 12 Phenanthrene, 3,6-dimethyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.30 | 15.44 ng | 938480 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 95 |
| 2 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 93 |
| 3 | Phenanthrene, 2,3-dimethyl- | 206 | C16H14 | 003674-65-5 | 93 |
| 4 | Phenanthrene, 2,7-dimethyl- | 206 | C16H14 | 001576-69-8 | 81 |
| 5 | di-p-Tolylacetylene | 206 | C16H14 | 002789-88-0 | 80 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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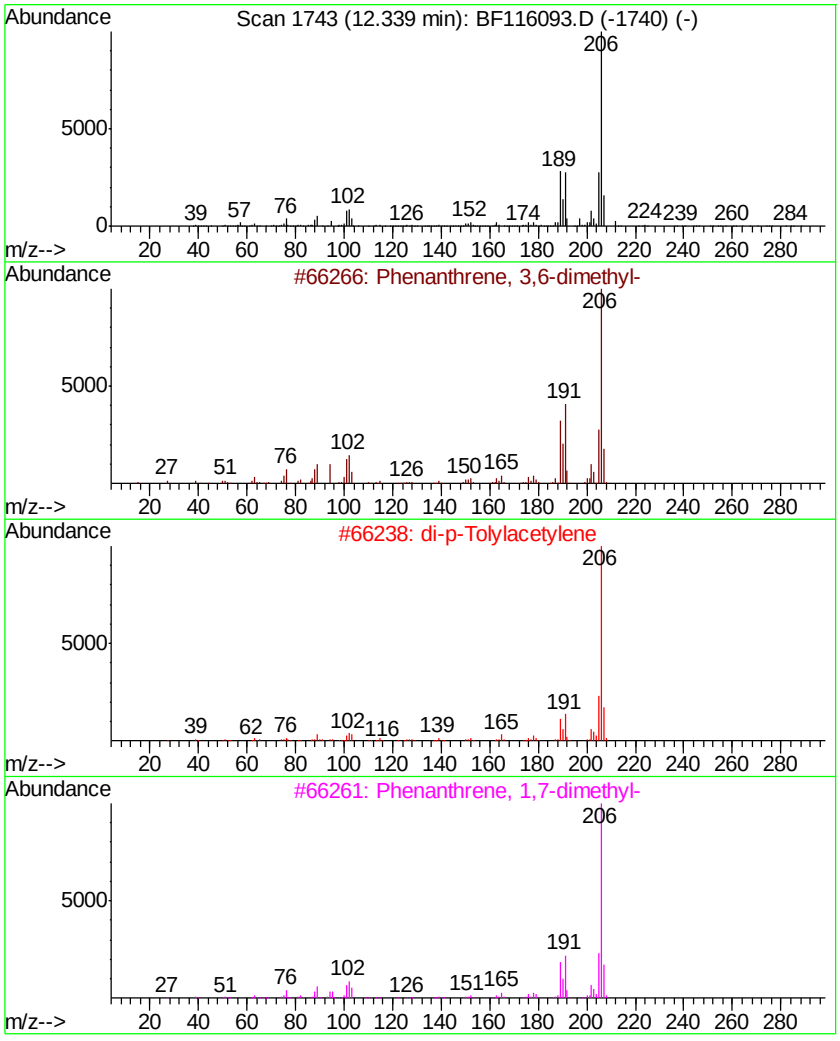
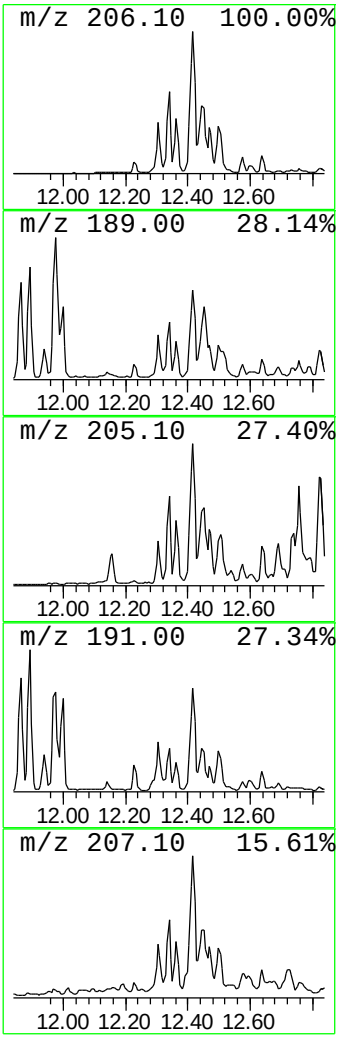
Peak Number 13 di-p-Tolylacetylene Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.34 | 13.28 ng | 807586 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 93 |
| 2 | di-p-Tolylacetylene | 206 | C16H14 | 002789-88-0 | 93 |
| 3 | Phenanthrene, 1,7-dimethyl- | 206 | C16H14 | 000483-87-4 | 93 |
| 4 | Phenanthrene, 2,7-dimethyl- | 206 | C16H14 | 001576-69-8 | 86 |
| 5 | Naphthalene, 1,2-dihydro-4-phenyl- | 206 | C16H14 | 007469-40-1 | 64 |



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Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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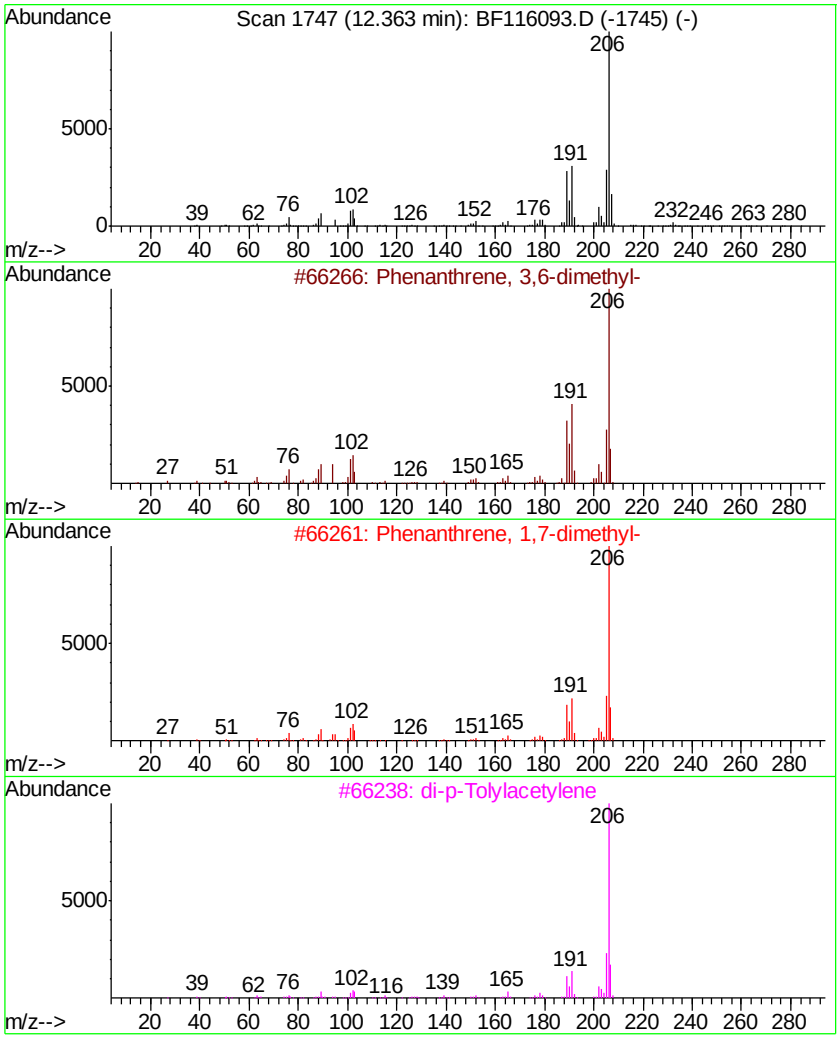
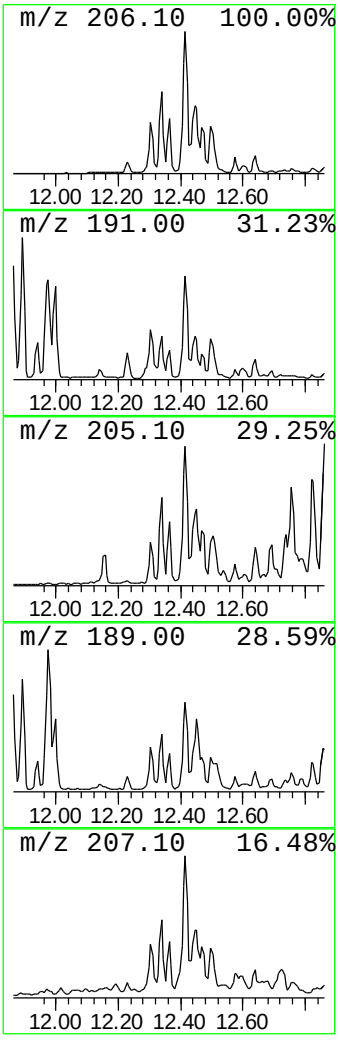
Peak Number 14 Phenanthrene, 1,7-dimethyl- Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.36 | 8.93 ng | 542808 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 95 |
| 2 | Phenanthrene, 1,7-dimethyl- | 206 | C16H14 | 000483-87-4 | 95 |
| 3 | di-p-Tolylacetylene | 206 | C16H14 | 002789-88-0 | 94 |
| 4 | Phenanthrene, 2,7-dimethyl- | 206 | C16H14 | 001576-69-8 | 93 |
| 5 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 90 |



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Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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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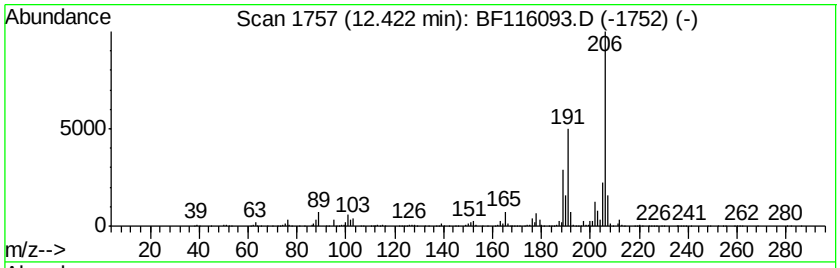
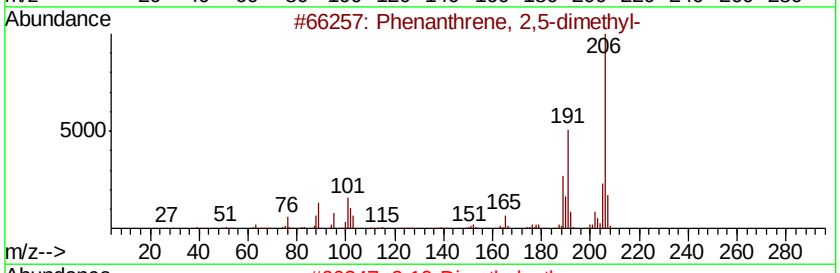
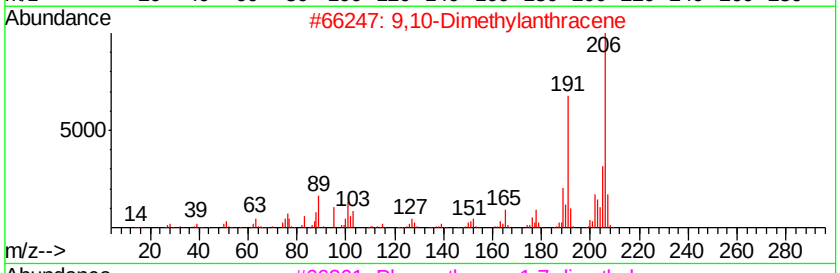
Peak Number 15 Phenanthrene, 2,5-dimethyl- Concentration Rank 1

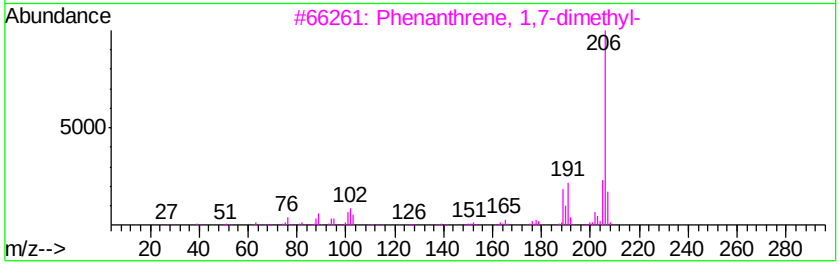
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.42 | 32.83 ng | 1996070 | Phenanthrene-d10 | 11.36 |

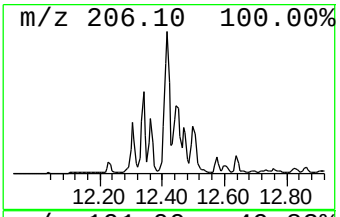
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 98 |
| 2 | 9,10-Dimethylanthracene | 206 | C16H14 | 000781-43-1 | 93 |
| 3 | Phenanthrene, 1,7-dimethyl- | 206 | C16H14 | 000483-87-4 | 90 |
| 4 | Phenanthrene, 2,3-dimethyl- | 206 | C16H14 | 003674-65-5 | 87 |
| 5 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 87 |

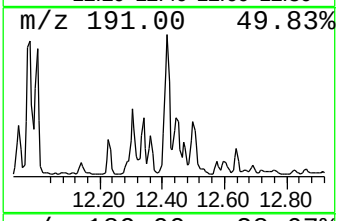


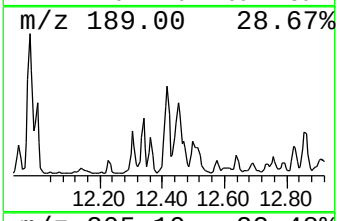



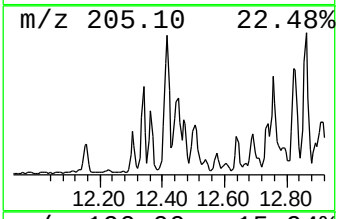



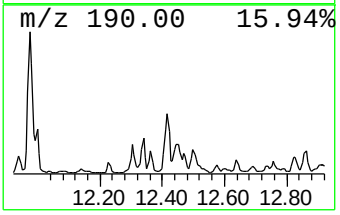



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

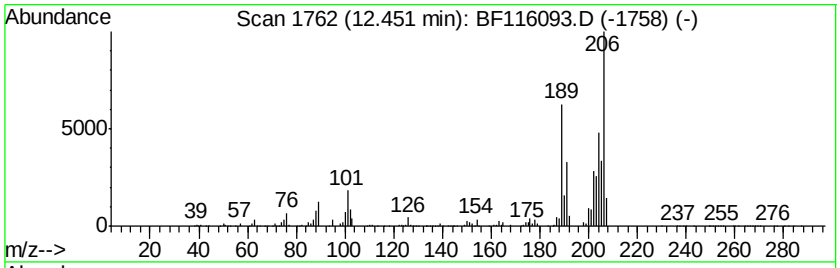
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

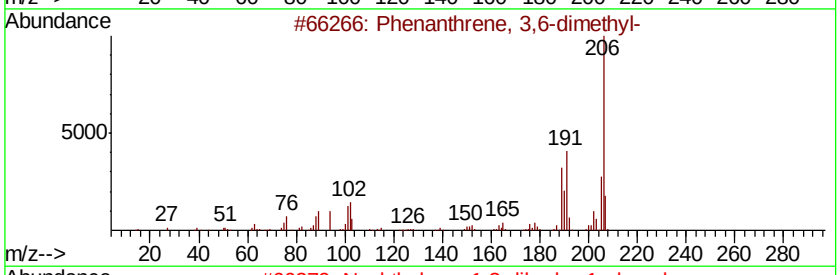
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

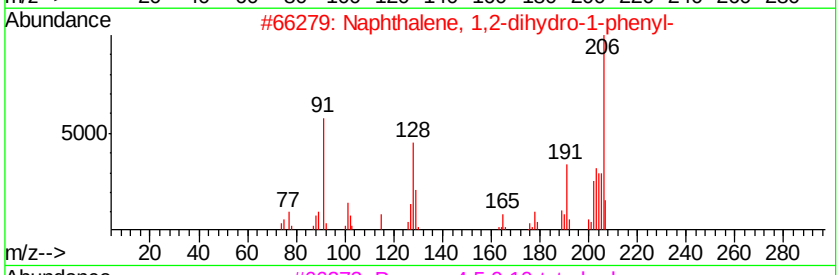
**********
Peak Number 16 Naphthalene, 1,2-dihydro-1-... Concentration Rank 8

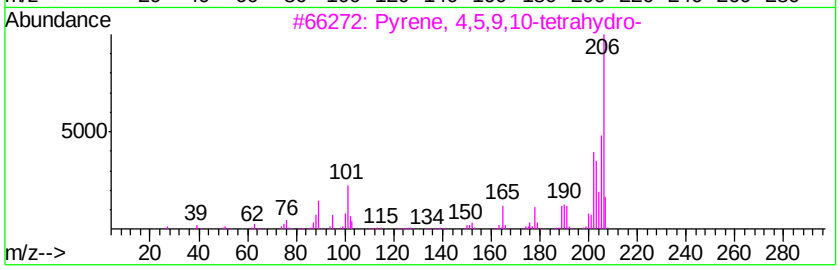
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.45 | 21.40 ng | 1301090 | Phenanthrene-d10 | 11.36 |

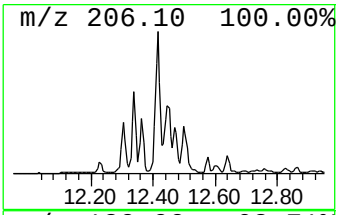
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 78 |
| 2 | Naphthalene, 1,2-dihydro-1-phenyl- | 206 | C16H14 | 016606-46-5 | 70 |
| 3 | Pyrene, 4,5,9,10-tetrahydro- | 206 | C16H14 | 000781-17-9 | 60 |
| 4 | 1,3-Butadiene, 1,4-diphenyl-, (E... | 206 | C16H14 | 000538-81-8 | 49 |
| 5 | Anthracene, 1,4-dimethyl- | 206 | C16H14 | 000781-92-0 | 46 |

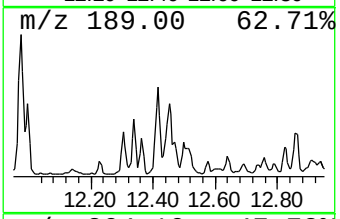


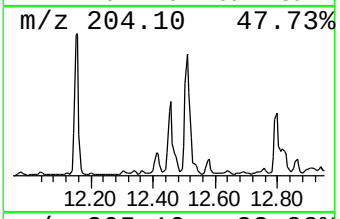



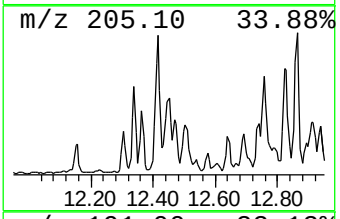



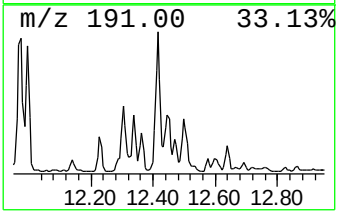



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

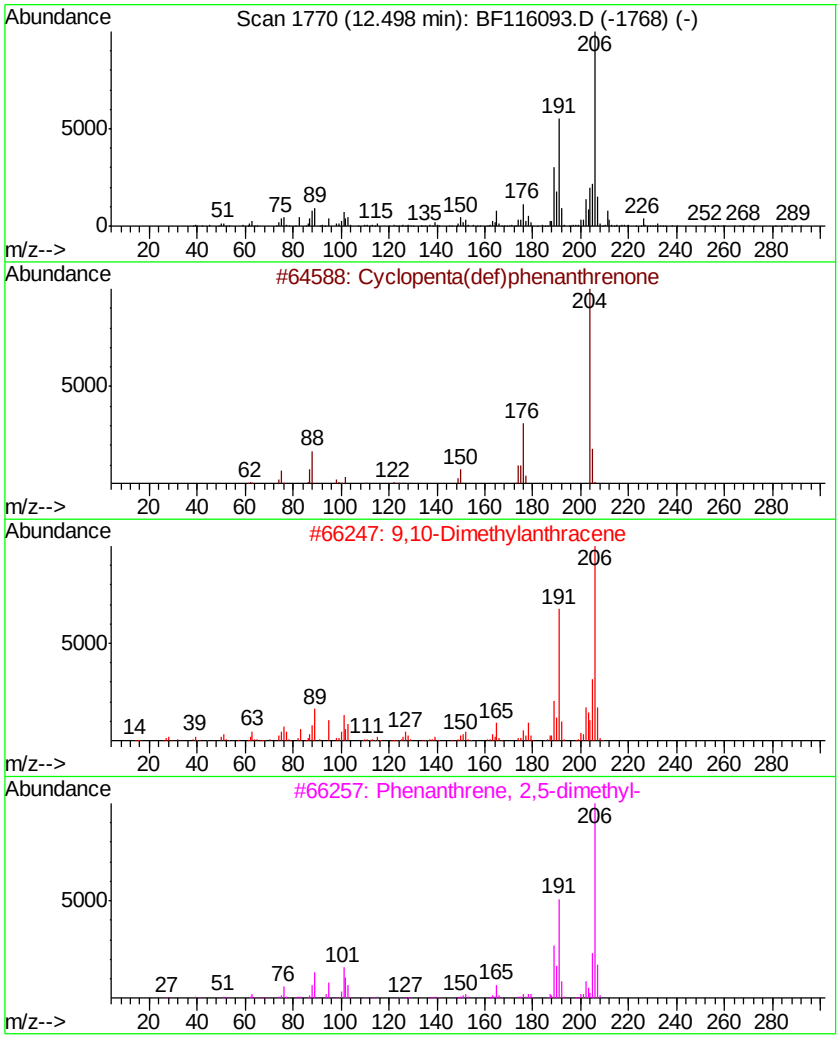
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

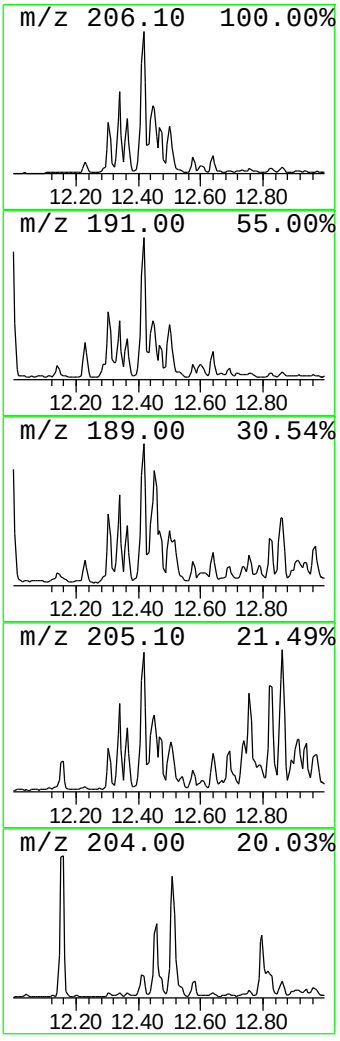
**********

Peak Number 17 Cyclopenta(def)phenanthrenone Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.50 | 16.31 ng | 991704 | Phenanthrene-d10 | 11.36 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopenta(def)phenanthrenone | 204 | C15H8O | 005737-13-3 | 74 |
| 2 | 9,10-Dimethylanthracene | 206 | C16H14 | 000781-43-1 | 55 |
| 3 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 53 |
| 4 | 1,3-Diphenyl-3-methylcyclopropene | 206 | C16H14 | 086544-79-8 | 49 |
| 5 | 1H-Indene, 2-methyl-3-phenyl- | 206 | C16H14 | 035099-60-6 | 49 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

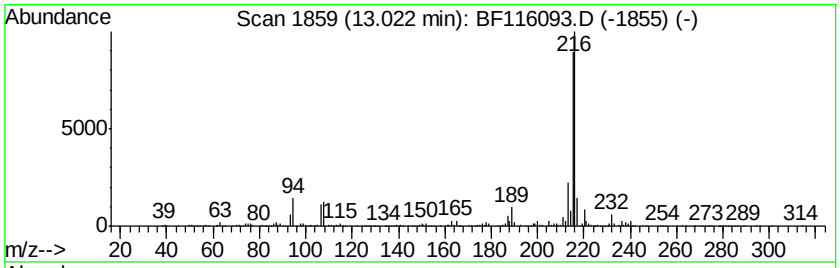
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

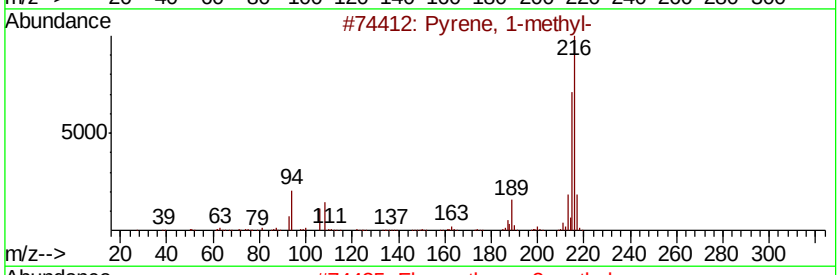
**********

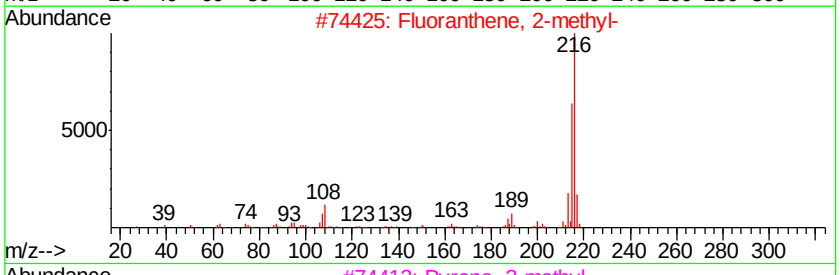
Peak Number 18 Fluoranthene, 2-methyl- Concentration Rank 16

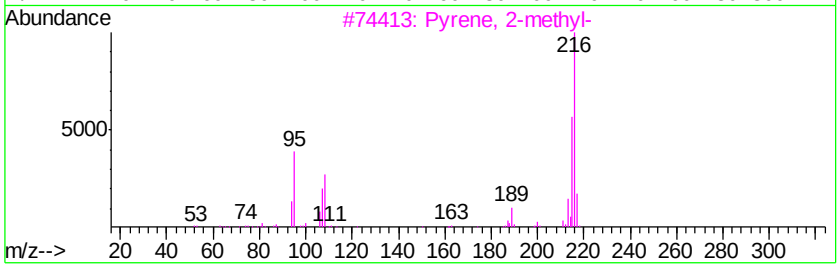
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.02 | 9.45 ng | 1415400 | Chrysene-d12 | 14.00 |

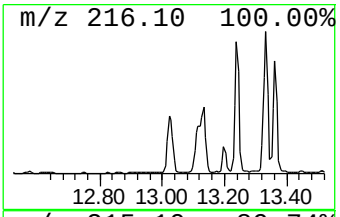
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 94 |
| 2 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 91 |
| 3 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 86 |
| 4 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 81 |
| 5 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 80 |

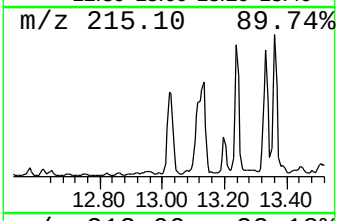


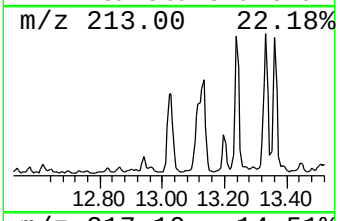



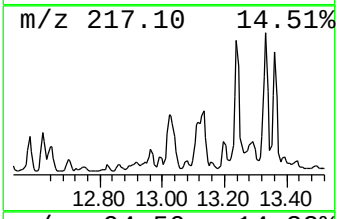



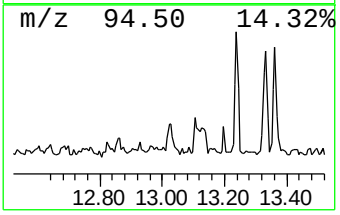



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

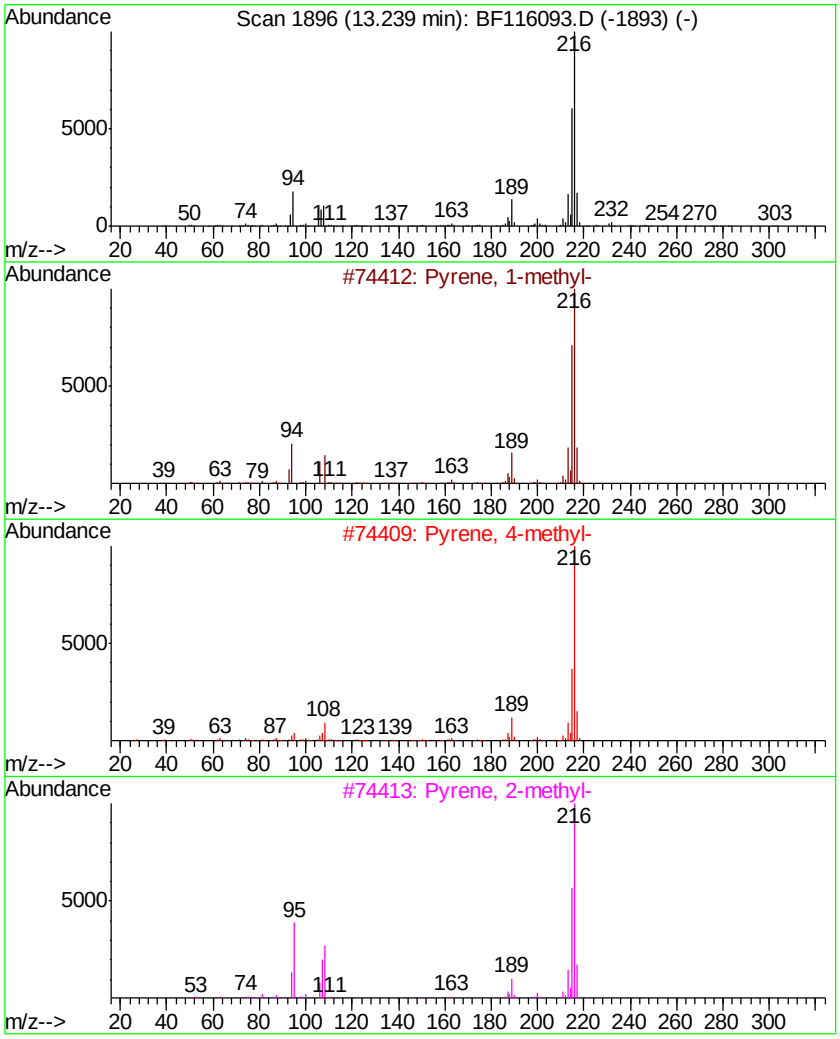
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

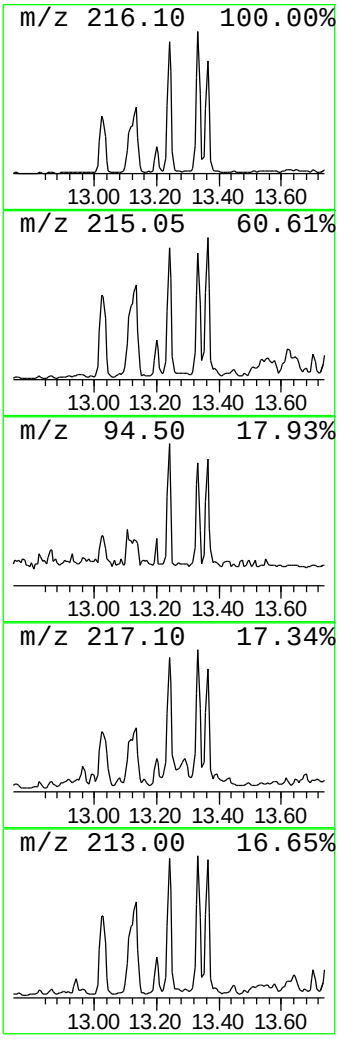
**********

Peak Number 19 Pyrene, 1-methyl- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.24 | 9.10 ng | 1363950 | Chrysene-d12 | 14.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 98 |
| 2 | Pyrene, 4-methyl- | 216 | C17H12 | 003353-12-6 | 94 |
| 3 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 94 |
| 4 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 90 |
| 5 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 90 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

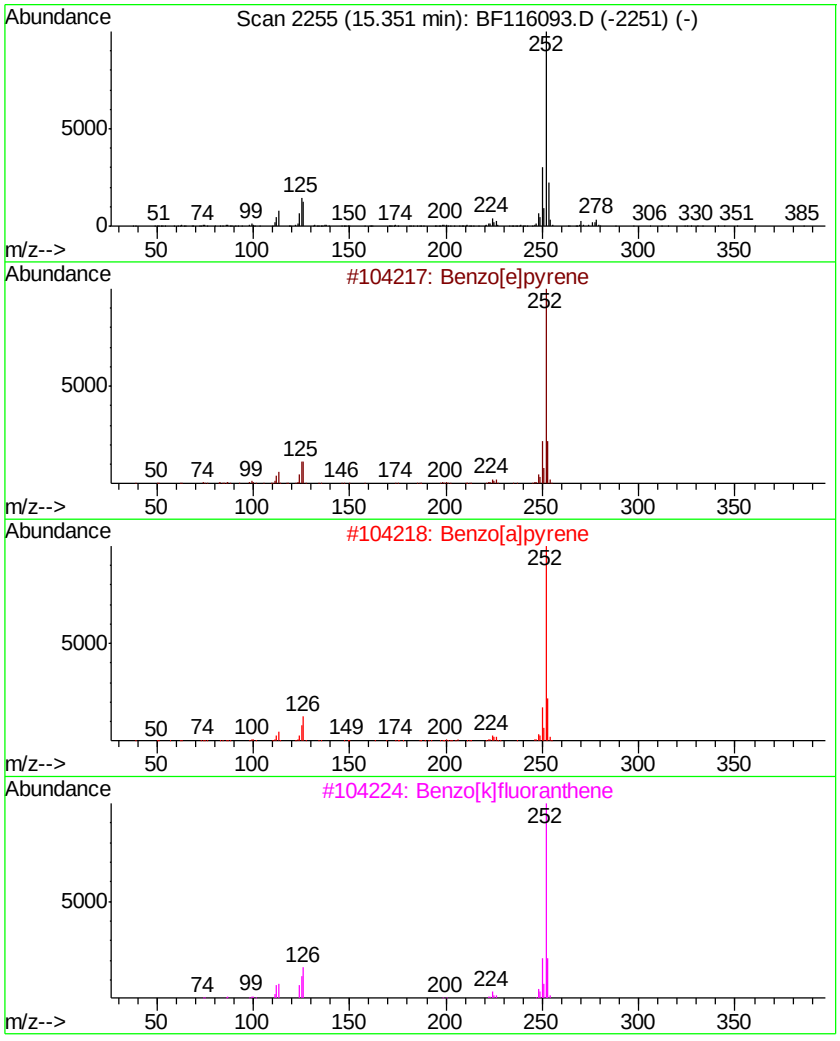
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

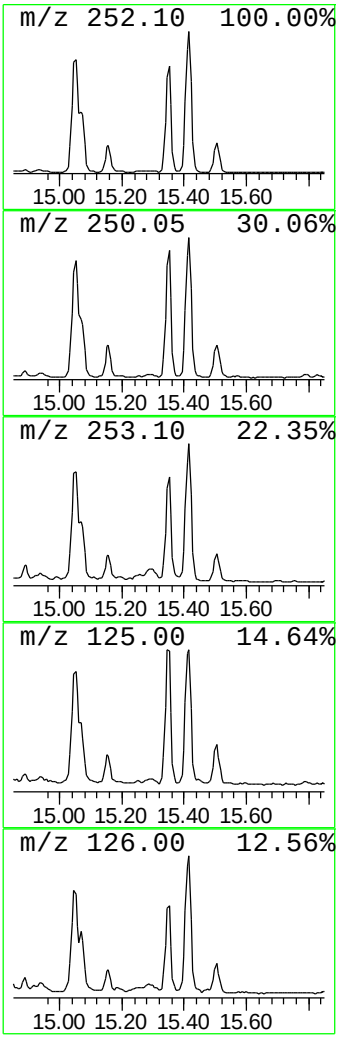
**********

Peak Number 20 Benzo[e]pyrene Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.35 | 24.82 ng | 1109700 | Perylene-d12 | 15.47 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 98 |
| 2 | Benzo[a]pyrene | 252 | C20H12 | 000050-32-8 | 98 |
| 3 | Benzo[k]fluoranthene | 252 | C20H12 | 000207-08-9 | 95 |
| 4 | Perylene | 252 | C20H12 | 000198-55-0 | 95 |
| 5 | Benz[e]acephenanthrylene | 252 | C20H12 | 000205-99-2 | 94 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080819\
Data File : BF116093.D
Acq On : 9 Aug 2019 2:58
Operator : HP/JU
Sample : K4178-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** 370-371

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| unknown6.61 | 6.61 | 32.0 | ng | 1453770 | 1 | 6.83 | 907676 | 20.0 |
| Naphthalene, 1,6,... | 10.19 | 8.6 | ng | 667442 | 3 | 9.87 | 1546810 | 20.0 |
| Chamazulene | 10.79 | 15.1 | ng | 919176 | 4 | 11.36 | 1216040 | 20.0 |
| 1-Methyldibenzoth... | 11.69 | 11.3 | ng | 685063 | 4 | 11.36 | 1216040 | 20.0 |
| Anthracene, 2-met... | 11.86 | 21.7 | ng | 1317760 | 4 | 11.36 | 1216040 | 20.0 |
| Phenanthrene, 2-m... | 11.89 | 27.1 | ng | 1647460 | 4 | 11.36 | 1216040 | 20.0 |
| Anthracene, 1-met... | 11.94 | 8.3 | ng | 505272 | 4 | 11.36 | 1216040 | 20.0 |
| Phenanthrene, 1-m... | 11.97 | 29.1 | ng | 1768740 | 4 | 11.36 | 1216040 | 20.0 |
| 1H-Indene, 2-phenyl- | 12.00 | 14.7 | ng | 891754 | 4 | 11.36 | 1216040 | 20.0 |
| Naphthalene, 2-ph... | 12.15 | 11.9 | ng | 724272 | 4 | 11.36 | 1216040 | 20.0 |
| Phenanthrene, 3,6... | 12.30 | 15.4 | ng | 938480 | 4 | 11.36 | 1216040 | 20.0 |
| di-p-Tolylacetylene | 12.34 | 13.3 | ng | 807586 | 4 | 11.36 | 1216040 | 20.0 |
| Phenanthrene, 1,7... | 12.36 | 8.9 | ng | 542808 | 4 | 11.36 | 1216040 | 20.0 |
| Phenanthrene, 2,5... | 12.42 | 32.8 | ng | 1996070 | 4 | 11.36 | 1216040 | 20.0 |
| Naphthalene, 1,2-... | 12.45 | 21.4 | ng | 1301090 | 4 | 11.36 | 1216040 | 20.0 |
| Cyclopenta(def)ph... | 12.50 | 16.3 | ng | 991704 | 4 | 11.36 | 1216040 | 20.0 |
| Fluoranthene, 2-m... | 13.02 | 9.4 | ng | 1415400 | 5 | 14.00 | 2996380 | 20.0 |
| Pyrene, 1-methyl- | 13.24 | 9.1 | ng | 1363950 | 5 | 14.00 | 2996380 | 20.0 |
| Benzo[e]pyrene | 15.35 | 24.8 | ng | 1109700 | 6 | 15.47 | 894031 | 20.0 |