LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

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-BF111319.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 | 2.158 | 8 | 12 | 18 | rVB | 444658 | 567286 | 25.14% | 1.617% |
| 2 | 5.093 | 507 | 511 | 518 | rVB | 243007 | 246901 | 10.94% | 0.704% |
| 3 | 5.498 | 576 | 580 | 584 | rBV | 1533238 | 1318954 | 58.46% | 3.759% |
| 4 | 6.510 | 748 | 752 | 762 | rBV | 1683717 | 1489095 | 66.00% | 4.244% |
| 5 | 6.657 | 773 | 777 | 780 | rBV | 1888711 | 1522998 | 67.50% | 4.341% |
| 6 | 6.893 | 813 | 817 | 820 | rBV | 1164917 | 979199 | 43.40% | 2.791% |
| 7 | 7.045 | 839 | 843 | 847 | rBV | 1535526 | 1225073 | 54.30% | 3.492% |
| 8 | 7.451 | 908 | 912 | 915 | rBV | 1084524 | 898271 | 39.81% | 2.560% |
| 9 | 8.175 | 1032 | 1035 | 1038 | rBV | 1339906 | 1219673 | 54.06% | 3.476% |
| 10 | 8.475 | 1083 | 1086 | 1089 | rVB2 | 28886 | 30544 | 1.35% | 0.087% |
| 11 | 8.892 | 1155 | 1157 | 1160 | rVB | 57257 | 40971 | 1.82% | 0.117% |
| 12 | 8.992 | 1170 | 1174 | 1177 | rBV | 52472 | 40960 | 1.82% | 0.117% |
| 13 | 9.257 | 1215 | 1219 | 1222 | rBV | 2104338 | 1778068 | 78.81% | 5.068% |
| 14 | 9.357 | 1234 | 1236 | 1240 | rVB | 28621 | 26337 | 1.17% | 0.075% |
| 15 | 9.516 | 1259 | 1263 | 1267 | rVB2 | 27822 | 31106 | 1.38% | 0.089% |
| 16 | 9.598 | 1274 | 1277 | 1279 | rBV | 30802 | 29908 | 1.33% | 0.085% |
| 17 | 9.645 | 1283 | 1285 | 1290 | rVB | 260146 | 217238 | 9.63% | 0.619% |
| 18 | 9.798 | 1308 | 1311 | 1315 | rBV | 30296 | 27502 | 1.22% | 0.078% |
| 19 | 9.939 | 1331 | 1335 | 1338 | rBV | 1561453 | 1273630 | 56.45% | 3.630% |
| 20 | 9.969 | 1338 | 1340 | 1344 | rVB | 189288 | 159790 | 7.08% | 0.455% |
| 21 | 10.145 | 1365 | 1370 | 1374 | rVB | 107396 | 99153 | 4.39% | 0.283% |
| 22 | 10.486 | 1425 | 1428 | 1434 | rBV | 198127 | 172302 | 7.64% | 0.491% |
| 23 | 10.645 | 1453 | 1455 | 1458 | rVB | 51481 | 42742 | 1.89% | 0.122% |
| 24 | 10.728 | 1465 | 1469 | 1473 | rBV | 1121720 | 930262 | 41.23% | 2.651% |
| 25 | 10.857 | 1480 | 1491 | 1494 | rBV5 | 35175 | 61013 | 2.70% | 0.174% |
| 26 | 11.039 | 1517 | 1522 | 1524 | rBV2 | 41299 | 47300 | 2.10% | 0.135% |
| 27 | 11.233 | 1552 | 1555 | 1560 | rVB | 76167 | 66565 | 2.95% | 0.190% |
| 28 | 11.328 | 1567 | 1571 | 1576 | rVB | 110995 | 103302 | 4.58% | 0.294% |
| 29 | 11.433 | 1585 | 1589 | 1591 | rBV | 1232834 | 1109562 | 49.18% | 3.162% |
| 30 | 11.457 | 1591 | 1593 | 1596 | rVV | 1721942 | 1329824 | 58.94% | 3.790% |
| 31 | 11.510 | 1596 | 1602 | 1605 | rVV | 463998 | 427972 | 18.97% | 1.220% |
| 32 | 11.539 | 1605 | 1607 | 1613 | rVB4 | 32643 | 31107 | 1.38% | 0.089% |
| 33 | 11.663 | 1622 | 1628 | 1631 | rBV2 | 165722 | 182749 | 8.10% | 0.521% |
| 34 | 11.727 | 1636 | 1639 | 1641 | rBV2 | 44880 | 39031 | 1.73% | 0.111% |

LSC Area Percent Report

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Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

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-BF111319.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 11.763 | 1643 | 1645 | 1652 | rVB3 | 35110 | 35212 | 1.56% | 0.100% |
| 36 | 11.822 | 1652 | 1655 | 1658 | rBV | 213521 | 170820 | 7.57% | 0.487% |
| 37 | 11.933 | 1669 | 1674 | 1676 | rBV | 153874 | 142687 | 6.32% | 0.407% |
| 38 | 11.963 | 1676 | 1679 | 1682 | rVV2 | 243450 | 293864 | 13.02% | 0.838% |
| 39 | 12.010 | 1684 | 1687 | 1690 | rVV | 75534 | 79010 | 3.50% | 0.225% |
| 40 | 12.051 | 1690 | 1694 | 1696 | rVV2 | 304824 | 298671 | 13.24% | 0.851% |
| 41 | 12.069 | 1696 | 1697 | 1701 | rVB | 105674 | 66934 | 2.97% | 0.191% |
| 42 | 12.133 | 1706 | 1708 | 1711 | rVB2 | 23185 | 23486 | 1.04% | 0.067% |
| 43 | 12.233 | 1722 | 1725 | 1727 | rBV | 104685 | 101683 | 4.51% | 0.290% |
| 44 | 12.257 | 1727 | 1729 | 1732 | rVB2 | 79992 | 70297 | 3.12% | 0.200% |
| 45 | 12.380 | 1745 | 1750 | 1753 | rVB3 | 54372 | 64078 | 2.84% | 0.183% |
| 46 | 12.410 | 1753 | 1755 | 1757 | rBV | 37483 | 35892 | 1.59% | 0.102% |
| 47 | 12.433 | 1757 | 1759 | 1763 | rVB2 | 29853 | 30898 | 1.37% | 0.088% |
| 48 | 12.486 | 1763 | 1768 | 1770 | rBV | 77753 | 96135 | 4.26% | 0.274% |
| 49 | 12.510 | 1770 | 1772 | 1773 | rVV | 253862 | 191762 | 8.50% | 0.547% |
| 50 | 12.527 | 1773 | 1775 | 1780 | rVV2 | 485724 | 494389 | 21.91% | 1.409% |
| 51 | 12.569 | 1780 | 1782 | 1787 | rVB2 | 87991 | 104542 | 4.63% | 0.298% |
| 52 | 12.616 | 1787 | 1790 | 1792 | rBV | 118308 | 111446 | 4.94% | 0.318% |
| 53 | 12.651 | 1792 | 1796 | 1800 | rBV | 2370398 | 2157649 | 95.63% | 6.149% |
| 54 | 12.739 | 1809 | 1811 | 1814 | rBV2 | 74473 | 76481 | 3.39% | 0.218% |
| 55 | 12.804 | 1819 | 1822 | 1825 | rVV | 90173 | 93847 | 4.16% | 0.267% |
| 56 | 12.880 | 1829 | 1835 | 1841 | rVV | 1976583 | 2256168 | 100.00% | 6.430% |
| 57 | 12.927 | 1841 | 1843 | 1847 | rVB2 | 82615 | 82869 | 3.67% | 0.236% |
| 58 | 13.022 | 1852 | 1859 | 1866 | rBV | 1525831 | 1491745 | 66.12% | 4.252% |
| 59 | 13.086 | 1867 | 1870 | 1871 | rBV | 64977 | 55083 | 2.44% | 0.157% |
| 60 | 13.127 | 1875 | 1877 | 1879 | rVB | 48317 | 33508 | 1.49% | 0.095% |
| 61 | 13.186 | 1883 | 1887 | 1888 | rBV3 | 58382 | 69176 | 3.07% | 0.197% |
| 62 | 13.204 | 1888 | 1890 | 1894 | rVB | 168279 | 166228 | 7.37% | 0.474% |
| 63 | 13.251 | 1894 | 1898 | 1899 | rBV3 | 24399 | 34607 | 1.53% | 0.099% |
| 64 | 13.274 | 1899 | 1902 | 1904 | rVB2 | 121154 | 92354 | 4.09% | 0.263% |
| 65 | 13.310 | 1904 | 1908 | 1911 | rBV2 | 108775 | 124726 | 5.53% | 0.355% |
| 66 | 13.339 | 1911 | 1913 | 1916 | rVB2 | 43447 | 40850 | 1.81% | 0.116% |
| 67 | 13.369 | 1916 | 1918 | 1921 | rBV2 | 25905 | 32497 | 1.44% | 0.093% |
| 68 | 13.404 | 1922 | 1924 | 1927 | rBV | 52882 | 42709 | 1.89% | 0.122% |
| 69 | 13.433 | 1927 | 1929 | 1933 | rBV3 | 54855 | 53741 | 2.38% | 0.153% |
| 70 | 13.716 | 1974 | 1977 | 1982 | rVB | 118025 | 116438 | 5.16% | 0.332% |
| 71 | 13.827 | 1992 | 1996 | 1999 | rBV2 | 144098 | 188885 | 8.37% | 0.538% |

LSC Area Percent Report

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Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

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-BF111319.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 13.857 | 1999 | 2001 | 2003 | rVV | 124067 | 109121 | 4.84% | 0.311% |
| 73 | 13.880 | 2003 | 2005 | 2009 | rVV2 | 125914 | 177628 | 7.87% | 0.506% |
| 74 | 13.927 | 2010 | 2013 | 2021 | rVB2 | 185561 | 255445 | 11.32% | 0.728% |
| 75 | 14.080 | 2031 | 2039 | 2042 | rBV3 | 1221117 | 1932787 | 85.67% | 5.509% |
| 76 | 14.110 | 2042 | 2044 | 2047 | rVB | 799401 | 588933 | 26.10% | 1.678% |
| 77 | 14.186 | 2055 | 2057 | 2061 | rBV | 149980 | 132663 | 5.88% | 0.378% |
| 78 | 14.304 | 2074 | 2077 | 2081 | rBV2 | 92495 | 115173 | 5.10% | 0.328% |
| 79 | 14.468 | 2103 | 2105 | 2108 | rVB | 112944 | 88832 | 3.94% | 0.253% |
| 80 | 14.504 | 2108 | 2111 | 2114 | rBV2 | 73809 | 70809 | 3.14% | 0.202% |
| 81 | 14.592 | 2124 | 2126 | 2128 | rBV | 83868 | 78648 | 3.49% | 0.224% |
| 82 | 15.139 | 2215 | 2219 | 2222 | rBV | 686297 | 1017770 | 45.11% | 2.901% |
| 83 | 15.210 | 2229 | 2231 | 2235 | rVV2 | 88935 | 82756 | 3.67% | 0.236% |
| 84 | 15.251 | 2236 | 2238 | 2241 | rVB | 153908 | 137152 | 6.08% | 0.391% |
| 85 | 15.457 | 2268 | 2273 | 2279 | rVB | 460311 | 648680 | 28.75% | 1.849% |
| 86 | 15.521 | 2279 | 2284 | 2290 | rBV | 668415 | 826546 | 36.63% | 2.356% |
| 87 | 15.586 | 2291 | 2295 | 2298 | rBV | 734235 | 899208 | 39.86% | 2.563% |
| 88 | 15.910 | 2348 | 2350 | 2353 | rBV3 | 85137 | 118951 | 5.27% | 0.339% |
| 89 | 17.104 | 2549 | 2553 | 2561 | rVB2 | 217478 | 418146 | 18.53% | 1.192% |

Sum of corrected areas: 35087003

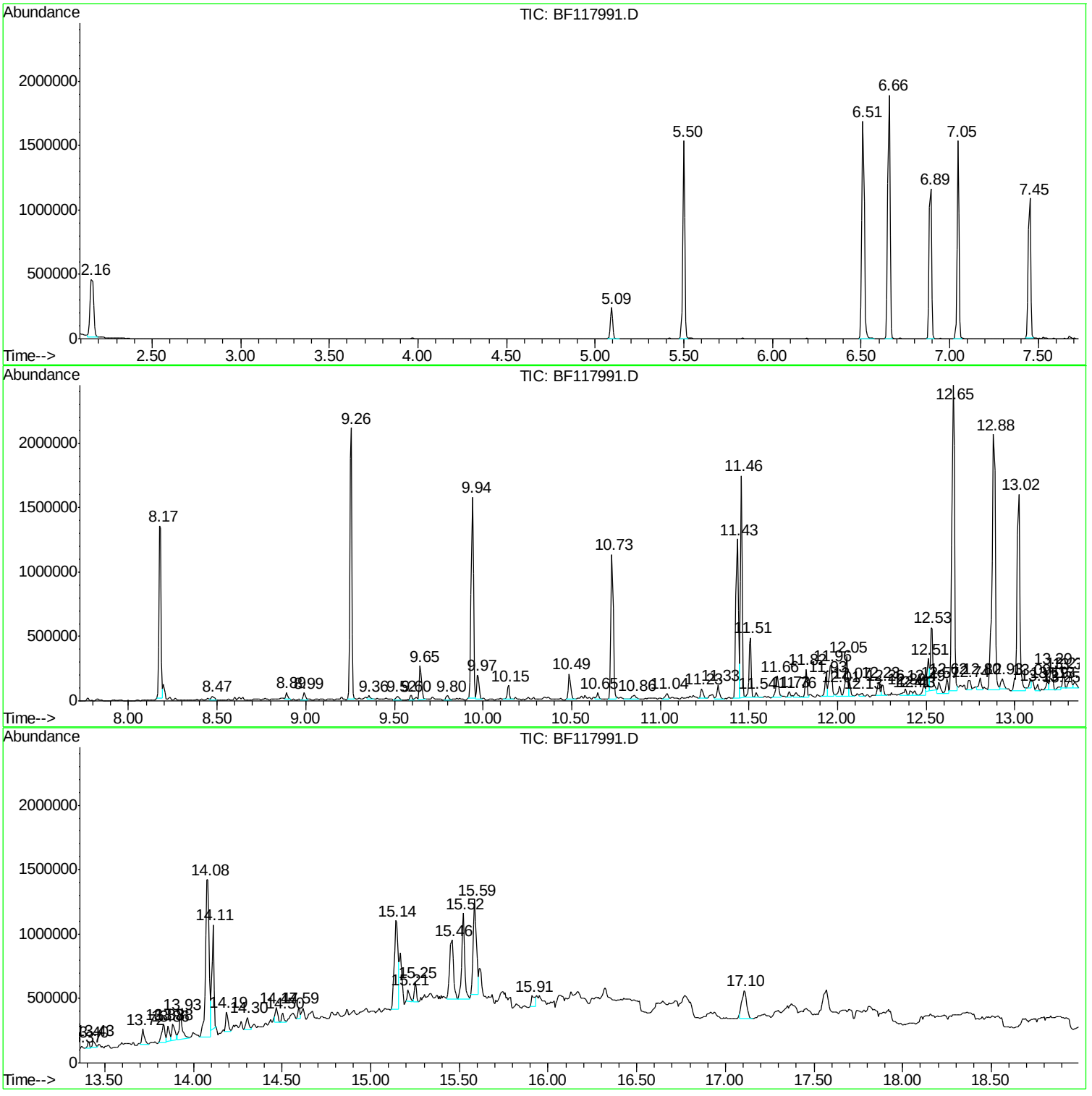
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.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\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

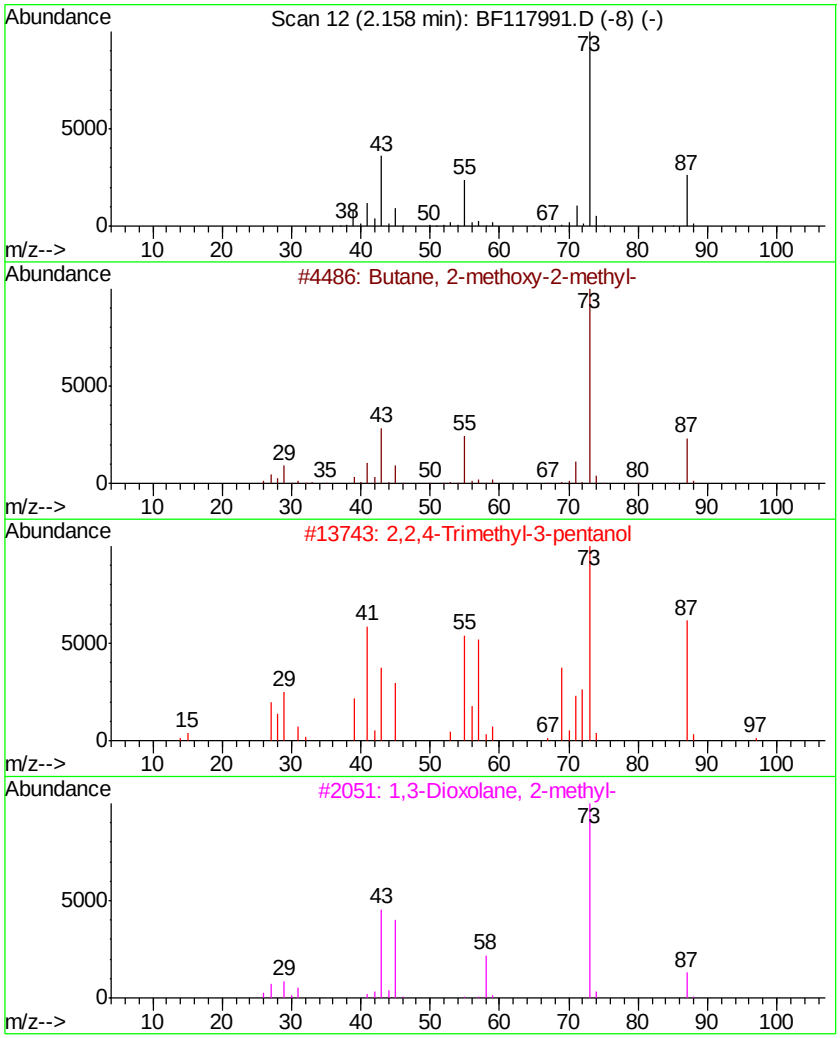
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

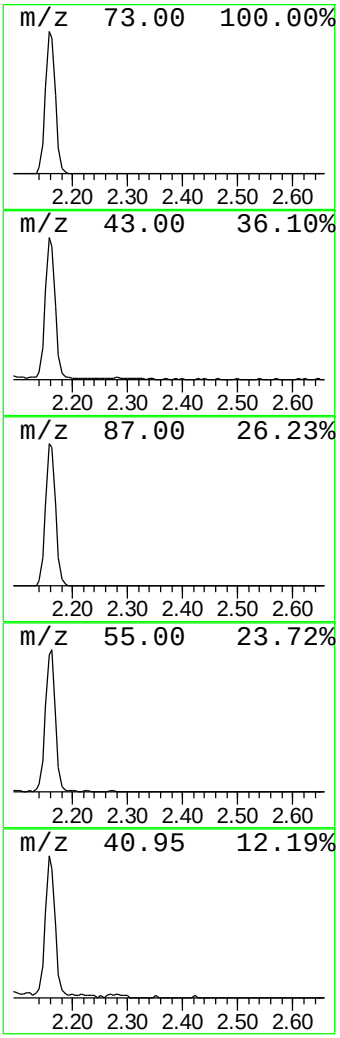
**********

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.16 | 11.59 ng | 567286 | 1,4-Dichlorobenzene-d4 | 6.89 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 78 |
| 2 | 2,2,4-Trimethyl-3-pentanol | 130 | C8H18O | 005162-48-1 | 56 |
| 3 | 1,3-Dioxolane, 2-methyl- | 88 | C4H8O2 | 000497-26-7 | 25 |
| 4 | Pentane, 3-methoxy- | 102 | C6H14O | 036839-67-5 | 23 |
| 5 | Silane, tetramethyl- | 88 | C4H12Si | 000075-76-3 | 17 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

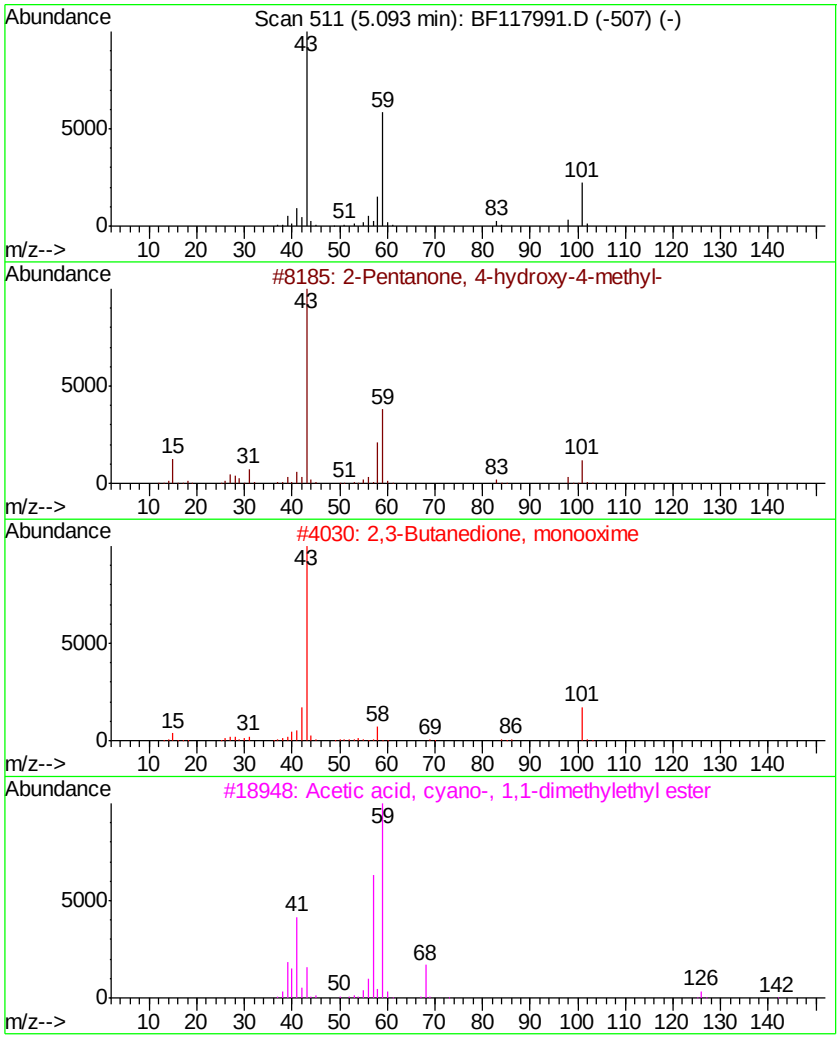
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

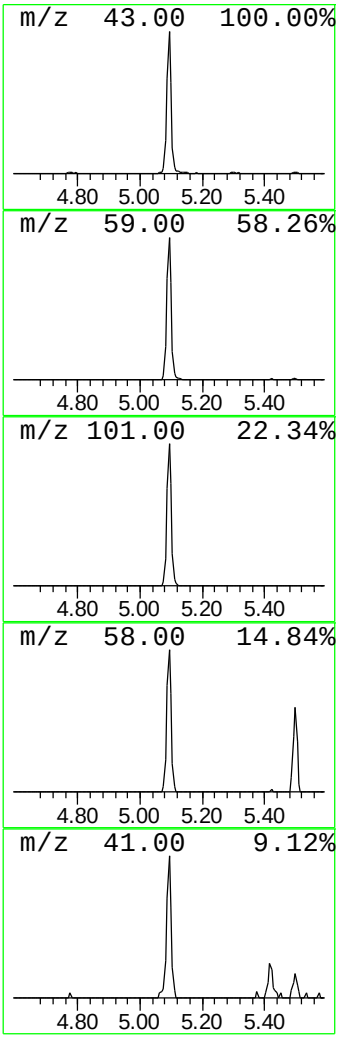
**********

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.09 | 5.04 ng | 246901 | 1,4-Dichlorobenzene-d4 | 6.89 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 56 |
| 2 | 2,3-Butanedione, monooxime | 101 | C4H7NO2 | 000057-71-6 | 35 |
| 3 | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 25 |
| 4 | Morpholine, 4-methyl- | 101 | C5H11NO | 000109-02-4 | 9 |
| 5 | 3-Hydroxy-3-methyl-2-butanone | 102 | C5H10O2 | 000115-22-0 | 9 |



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Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

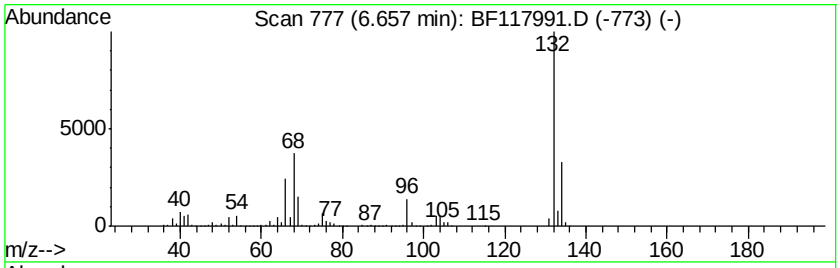
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

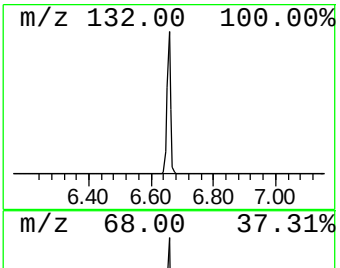
**********

Peak Number 3 unknown6.66 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.66 | 31.11 ng | 1523000 | 1,4-Dichlorobenzene-d4 | 6.89 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO | 031357-76-3 | 25 |
| 2 | Tranylcypromine-propionyl | 189 | C12H15NO | 1000123-86-3 | 16 |
| 3 | Cyclopropanamine, 1-phenyl- | 133 | C9H11N | 1000351-26-2 | 9 |
| 4 | Xenon | 132 | Xe | 007440-63-3 | 9 |
| 5 | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2 | 022291-85-6 | 9 |



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Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

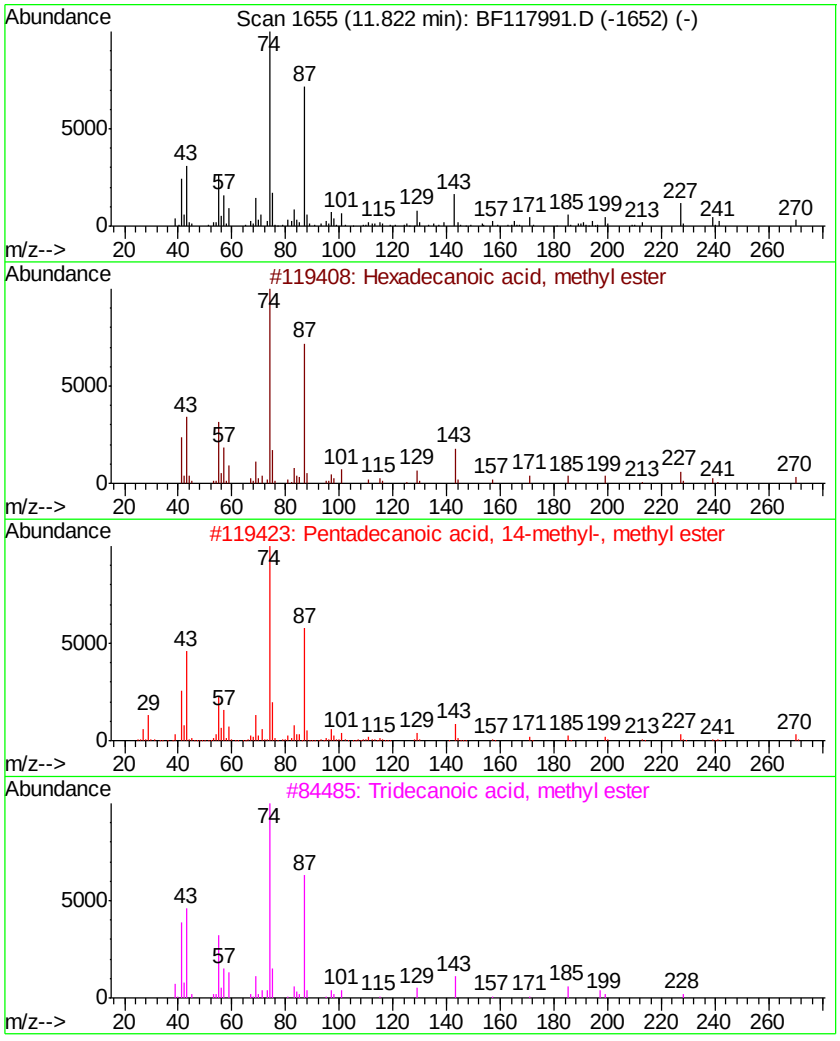
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

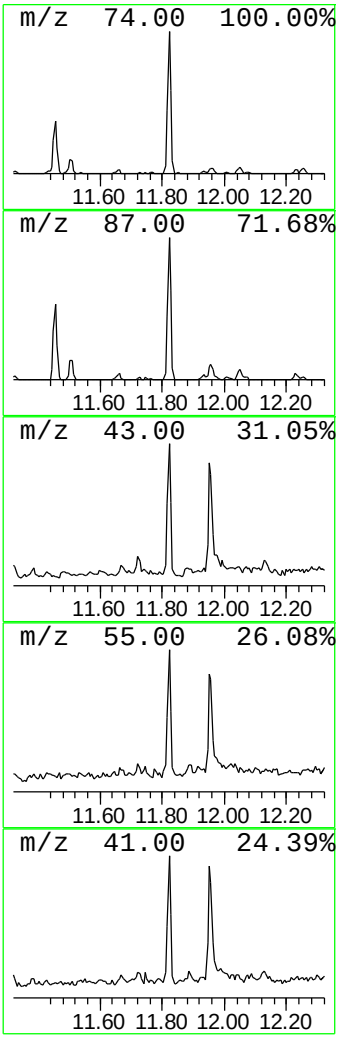
**********

Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.82 | 3.08 ng | 170820 | Phenanthrene-d10 | 11.43 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexadecanoic acid, methyl ester | 270 | C17H34O2 | 000112-39-0 | 98 |
| 2 | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 95 |
| 3 | Tridecanoic acid, methyl ester | 228 | C14H28O2 | 001731-88-0 | 91 |
| 4 | Pentadecanoic acid, methyl ester | 256 | C16H32O2 | 007132-64-1 | 80 |
| 5 | Dodecanoic acid, methyl ester | 214 | C13H26O2 | 000111-82-0 | 80 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

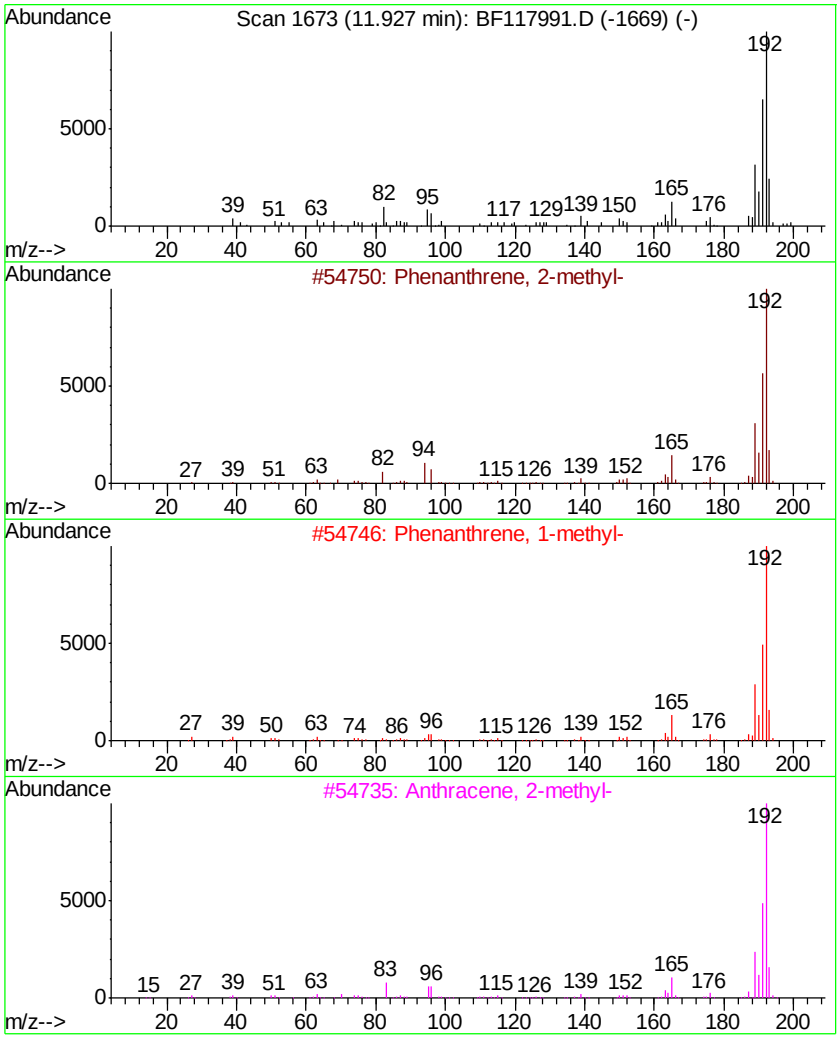
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

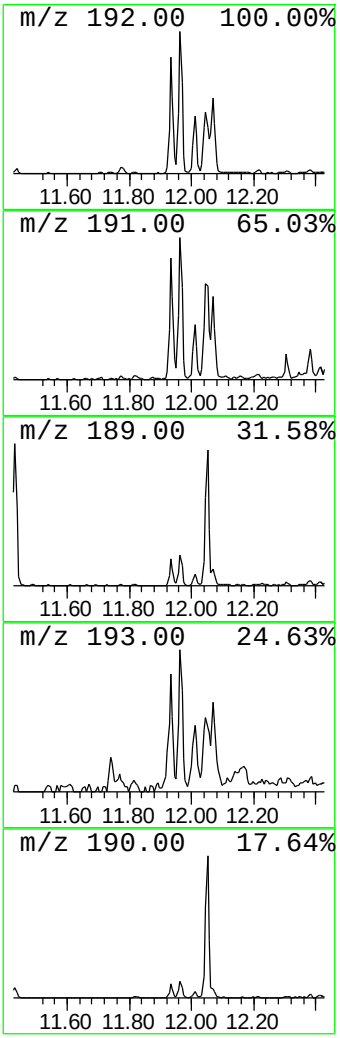
**********

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.93 | 2.57 ng | 142687 | Phenanthrene-d10 | 11.43 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 98 |
| 2 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 97 |
| 3 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 96 |
| 4 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 95 |
| 5 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 95 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

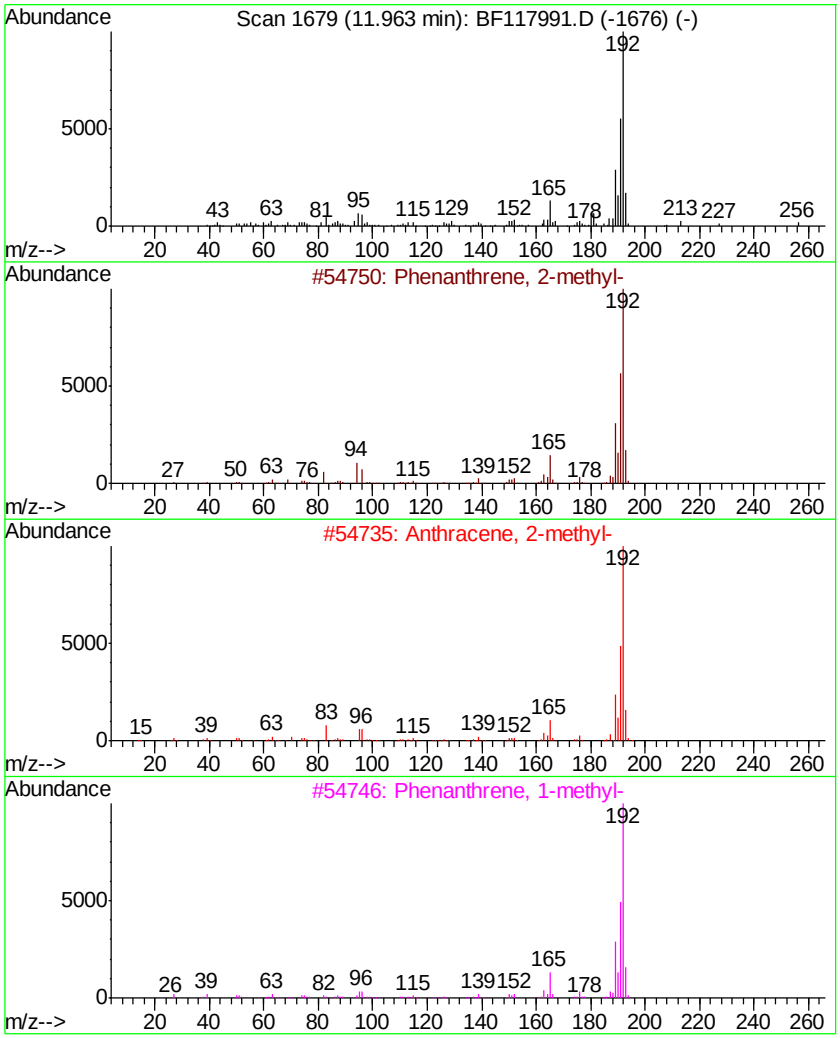
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

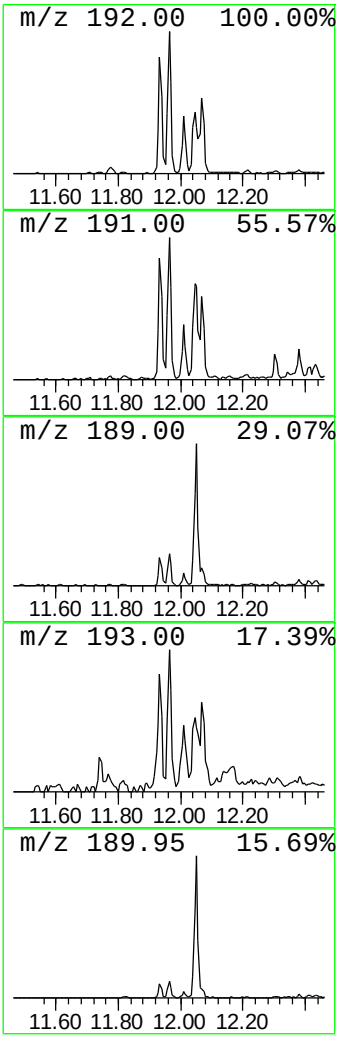
**********

Peak Number 7 Anthracene, 2-methyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.96 | 5.30 ng | 293864 | Phenanthrene-d10 | 11.43 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 97 |
| 2 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 96 |
| 3 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 95 |
| 4 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 93 |
| 5 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 91 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

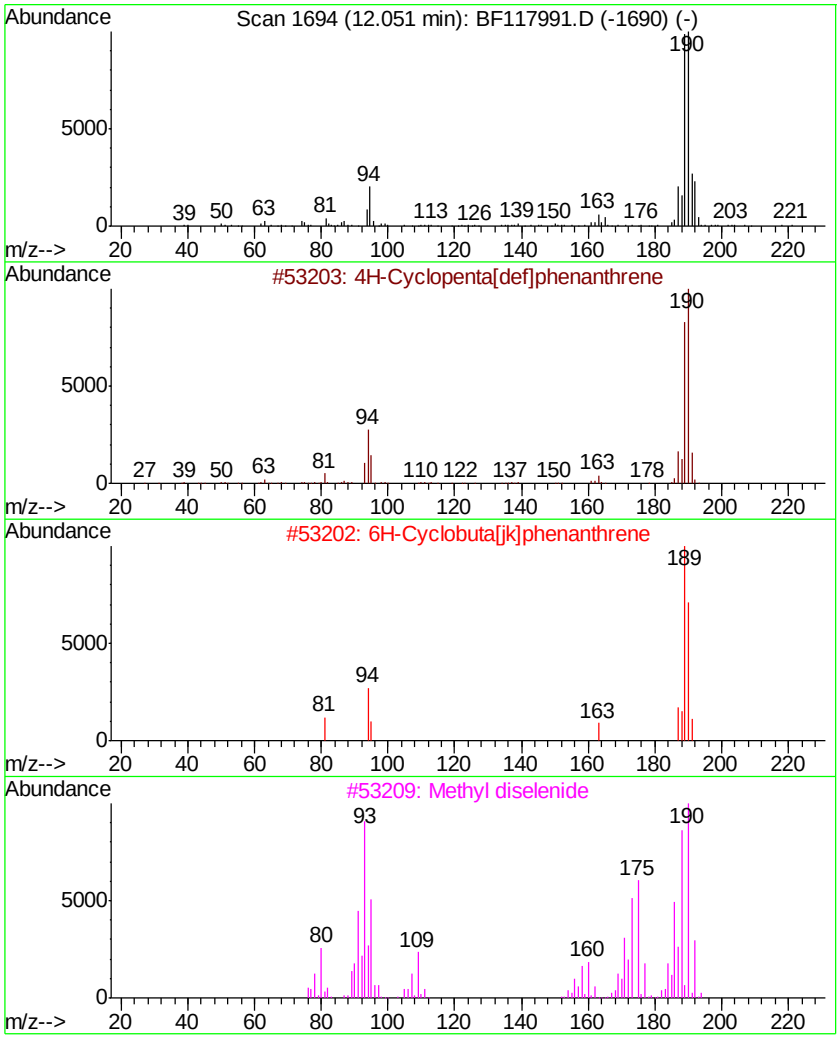
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

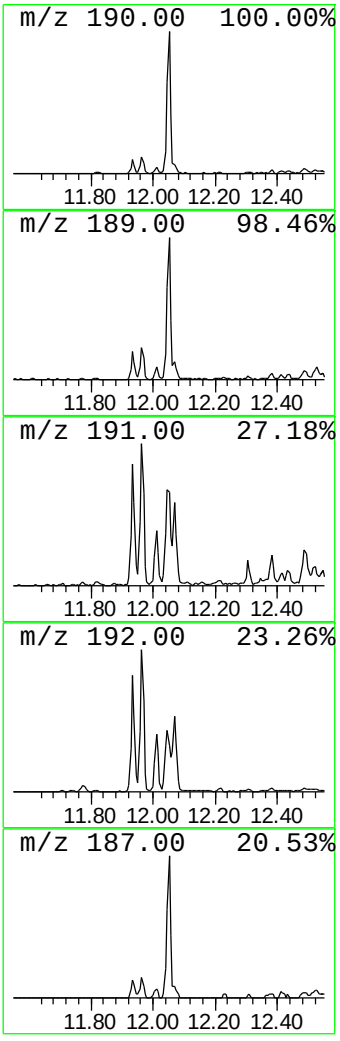
**********

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.05 | 5.38 ng | 298671 | Phenanthrene-d10 | 11.43 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 93 |
| 2 | 6H-Cyclobuta[jk]phenanthrene | 190 | C15H10 | 083469-43-6 | 64 |
| 3 | Methyl diselenide | 190 | C2H6Se2 | 007101-31-7 | 49 |
| 4 | 2,2'-Bis(4,5-dimethylimidazole) | 190 | C10H14N4 | 069286-06-2 | 42 |
| 5 | Megastigmatrienone | 190 | C13H18O | 038818-55-2 | 18 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

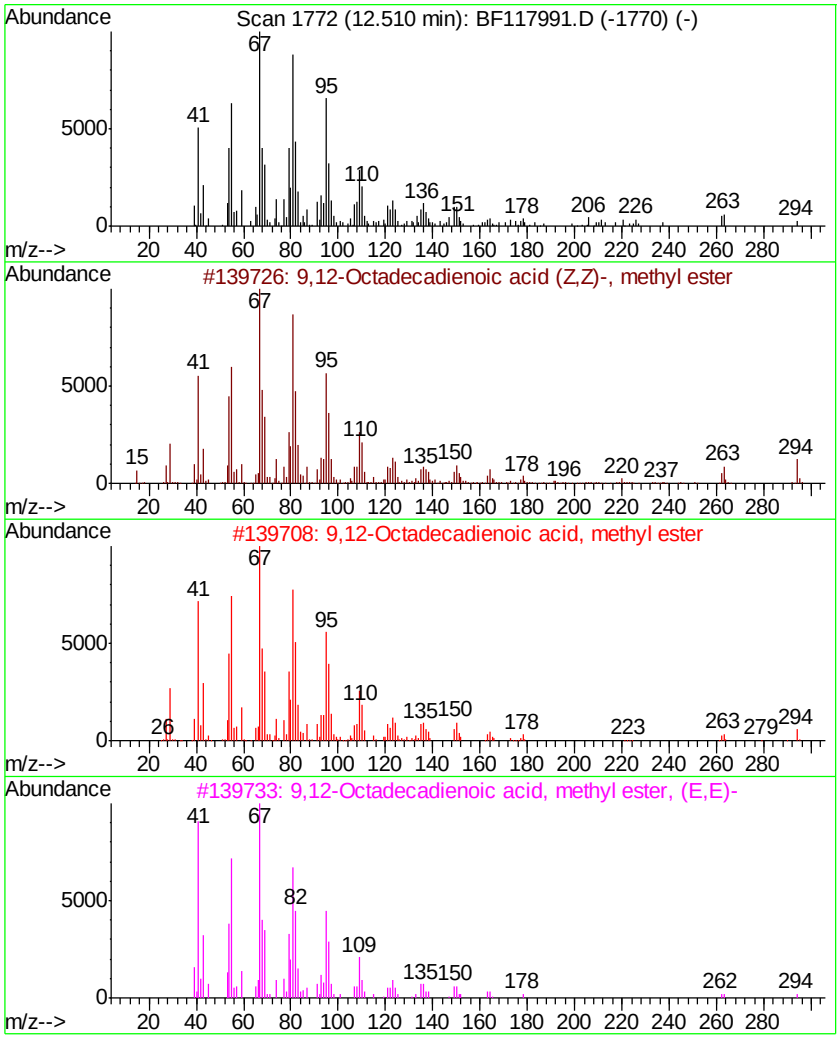
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

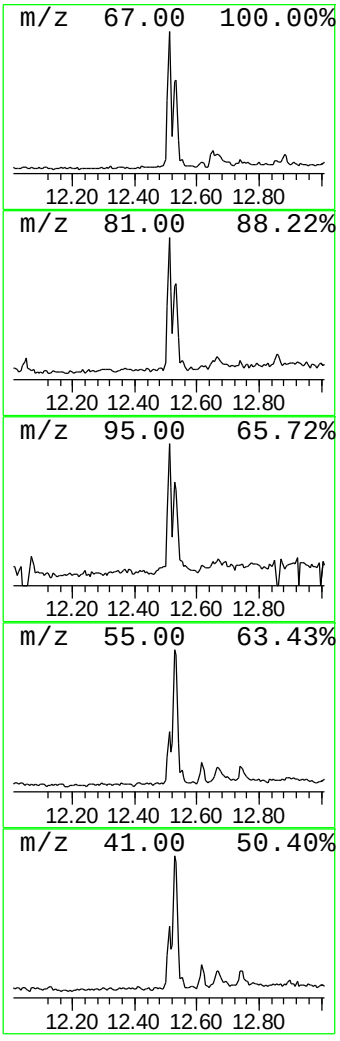
**********

Peak Number 9 9,12-Octadecadienoic acid (... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.51 | 3.46 ng | 191762 | Phenanthrene-d10 | 11.43 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9,12-Octadecadienoic acid (Z,Z)-... | 294 | C19H34O2 | 000112-63-0 | 99 |
| 2 | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002462-85-3 | 99 |
| 3 | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002566-97-4 | 99 |
| 4 | 11,14-Octadecadienoic acid, meth... | 294 | C19H34O2 | 056554-61-1 | 93 |
| 5 | 12,15-Octadecadienoic acid, meth... | 294 | C19H34O2 | 057156-97-5 | 91 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

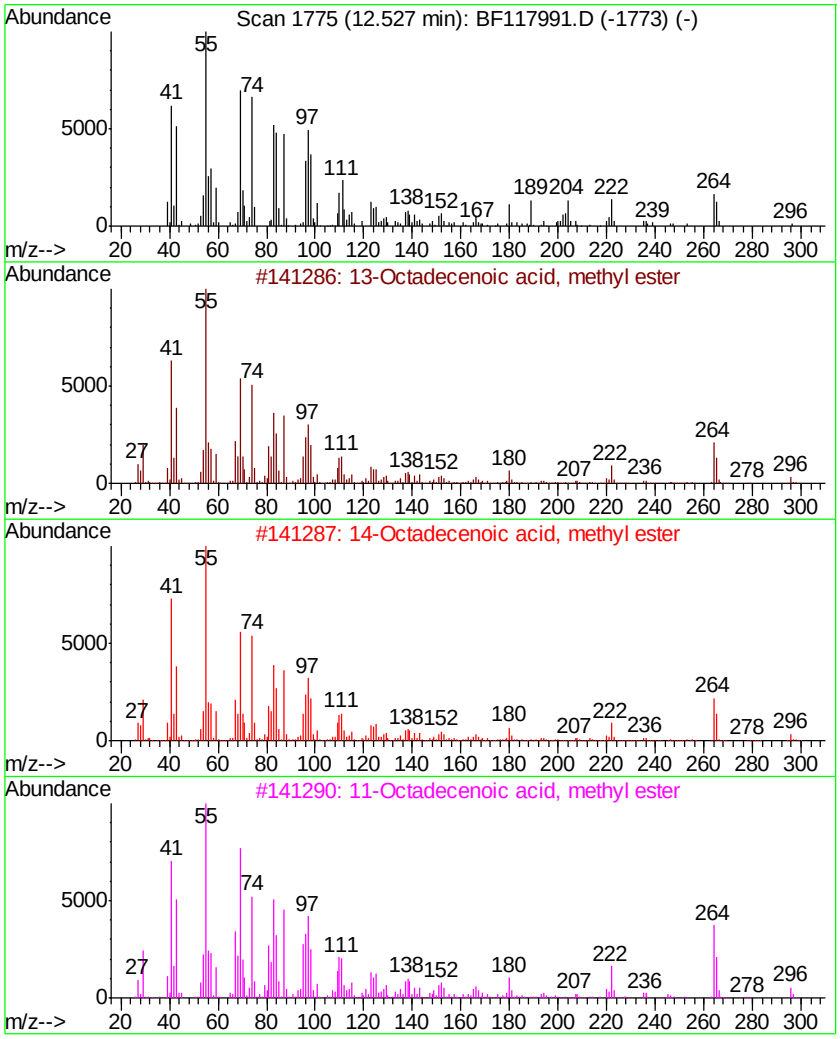
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

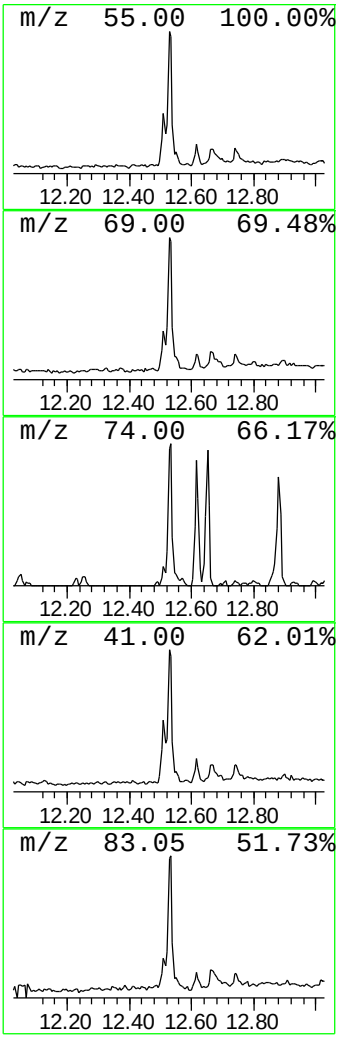
**********

Peak Number 10 13-Octadecenoic acid, methy... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.53 | 8.91 ng | 494389 | Phenanthrene-d10 | 11.43 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 13-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 056554-47-3 | 99 |
| 2 | 14-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 056554-48-4 | 97 |
| 3 | 11-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 052380-33-3 | 96 |
| 4 | 16-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 056554-49-5 | 95 |
| 5 | 15-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 004764-72-1 | 95 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

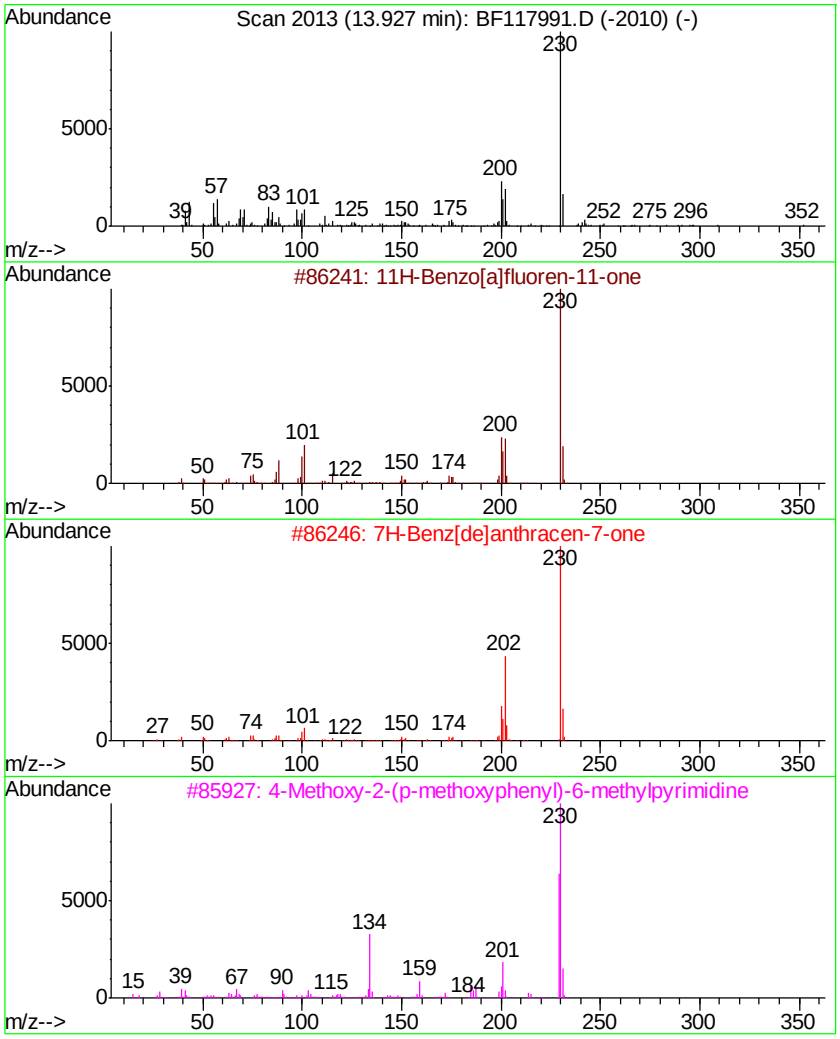
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

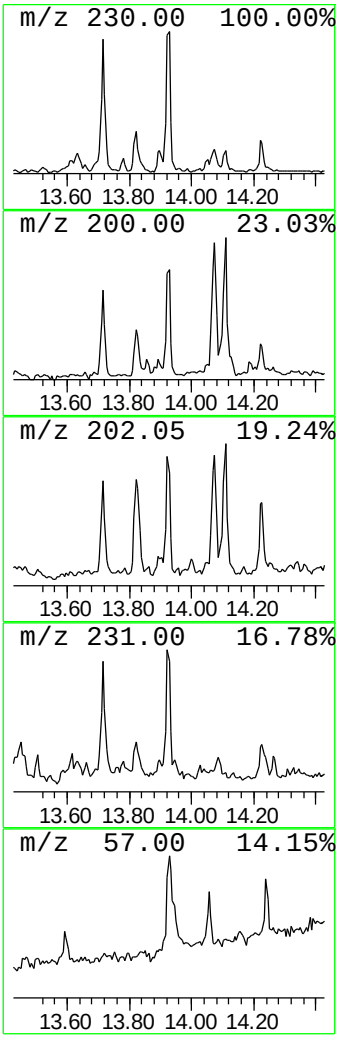
**********

Peak Number 11 11H-Benzo[a]fluoren-11-one Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.93 | 2.64 ng | 255445 | Chrysene-d12 | 14.09 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 11H-Benzo[a]fluoren-11-one | 230 | C17H10O | 000479-79-8 | 95 |
| 2 | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 89 |
| 3 | 4-Methoxy-2-(p-methoxyphenyl)-6-... | 230 | C13H14N2O2 | 063896-93-5 | 50 |
| 4 | Benzo(b)phenazine | 230 | C16H10N2 | 000257-97-6 | 47 |
| 5 | o-Terphenyl | 230 | C18H14 | 000084-15-1 | 47 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

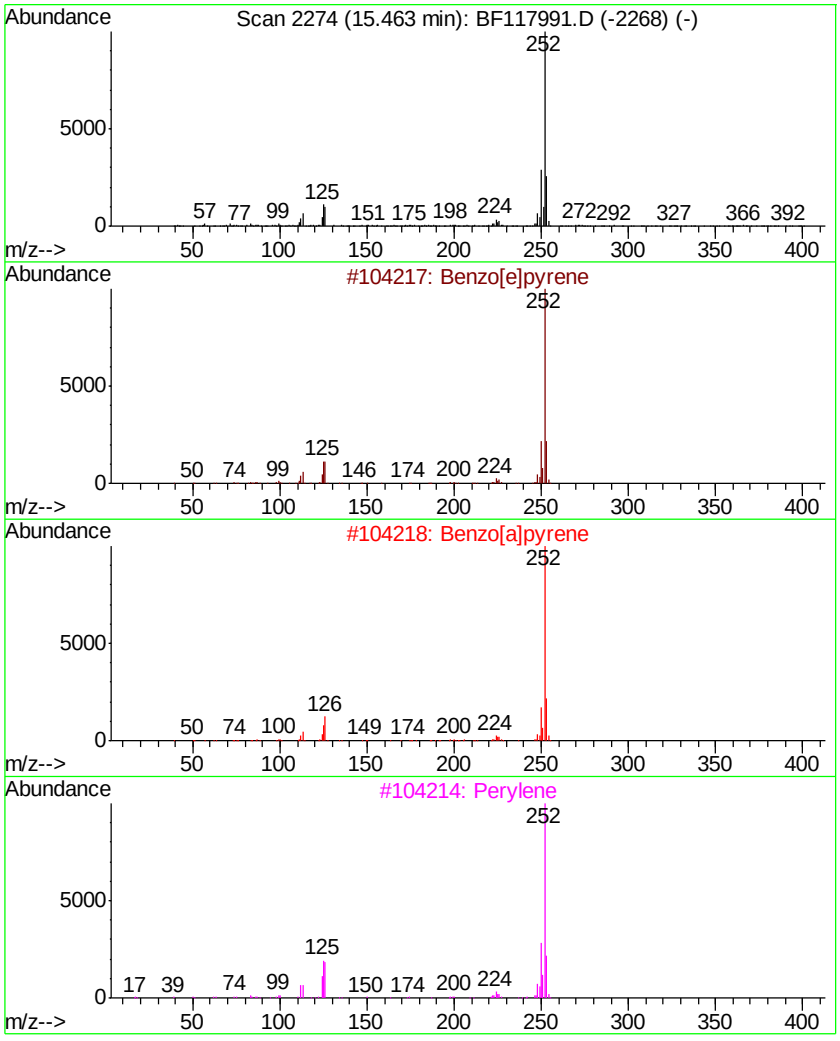
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

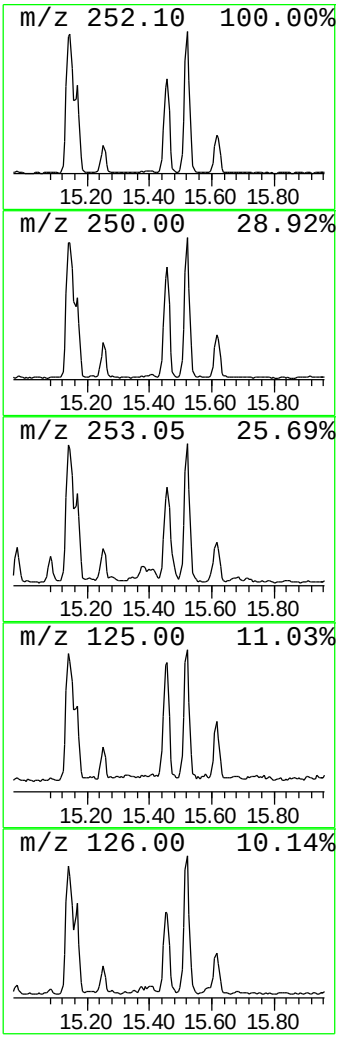
**********

Peak Number 13 Benzo[e]pyrene Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.46 | 14.43 ng | 648680 | Perylene-d12 | 15.59 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 99 |
| 2 | Benzo[a]pyrene | 252 | C20H12 | 000050-32-8 | 97 |
| 3 | Perylene | 252 | C20H12 | 000198-55-0 | 97 |
| 4 | Benz[e]acephenanthrylene | 252 | C20H12 | 000205-99-2 | 93 |
| 5 | Benzo[k]fluoranthene | 252 | C20H12 | 000207-08-9 | 93 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

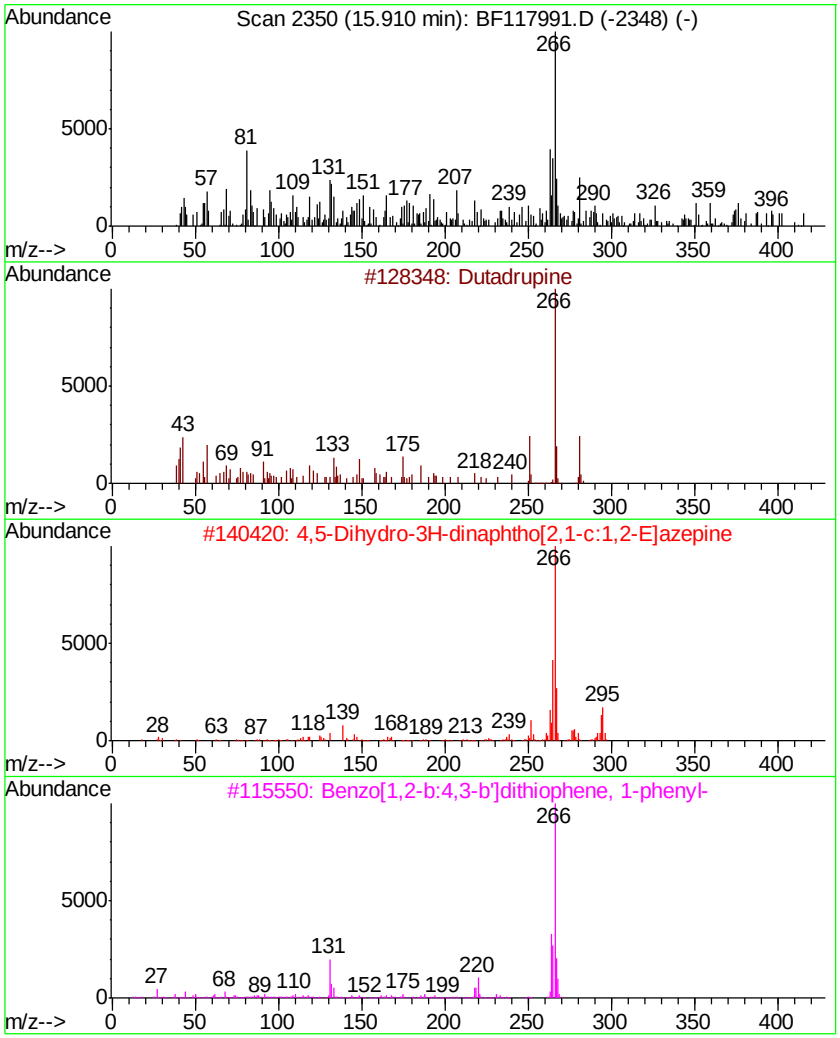
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

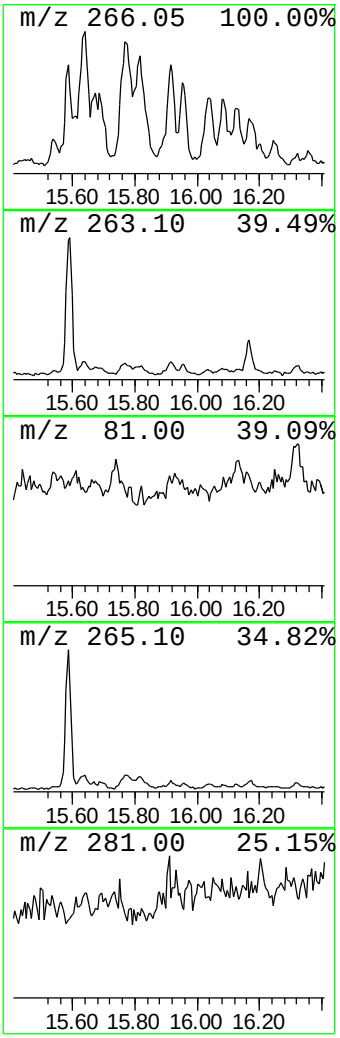
**********

Peak Number 14 Dutadrupine Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.91 | 2.65 ng | 118951 | Perylene-d12 | 15.59 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dutadrupine | 281 | C17H15NO3 | 080151-78-6 | 64 |
| 2 | 4,5-Dihydro-3H-dinaphtho[2,1-c:1... | 295 | C22H17N | 1000297-99-2 | 55 |
| 3 | Benzo[1,2-b:4,3-b']dithiophene, ... | 266 | C16H10S2 | 016587-58-9 | 50 |
| 4 | Benz[j]aceanthrylene, 3-methyl- | 266 | C21H14 | 003343-10-0 | 45 |
| 5 | 4'-Amino-4-dimethylaminochalcone | 266 | C17H18N2O | 025870-72-8 | 43 |



# Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF112519\
Data File : BF117991.D
Acq On : 26 Nov 2019 3:08
Operator : JU
Sample : K5663-01 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** AU-06-112219

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.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 |
| Butane, 2-methoxy... | 2.16 | 11.6 | ng | 567286 | 1 | 6.89 | 979199 | 20.0 |
| 2-Pentanone, 4-hy... | 5.09 | 5.0 | ng | 246901 | 1 | 6.89 | 979199 | 20.0 |
| unknown6.66 | 6.66 | 31.1 | ng | 1523000 | 1 | 6.89 | 979199 | 20.0 |
| Hexadecanoic acid... | 11.82 | 3.1 | ng | 170820 | 4 | 11.43 | 1109560 | 20.0 |
| Phenanthrene, 2-m... | 11.93 | 2.6 | ng | 142687 | 4 | 11.43 | 1109560 | 20.0 |
| Anthracene, 2-met... | 11.96 | 5.3 | ng | 293864 | 4 | 11.43 | 1109560 | 20.0 |
| 4H-Cyclopenta[def... | 12.05 | 5.4 | ng | 298671 | 4 | 11.43 | 1109560 | 20.0 |
| 9,12-Octadecadien... | 12.51 | 3.5 | ng | 191762 | 4 | 11.43 | 1109560 | 20.0 |
| 13-Octadecenoic a... | 12.53 | 8.9 | ng | 494389 | 4 | 11.43 | 1109560 | 20.0 |
| 11H-Benzo[a]fluor... | 13.93 | 2.6 | ng | 255445 | 5 | 14.09 | 1932790 | 20.0 |
| Benzo[e]pyrene | 15.46 | 14.4 | ng | 648680 | 6 | 15.59 | 899208 | 20.0 |
| Dutadrupine | 15.91 | 2.6 | ng | 118951 | 6 | 15.59 | 899208 | 20.0 |