LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

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-BF083019.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 | 5.451 | 564 | 572 | 578 | rBV | 113995 | 106858 | 4.67% | 0.346% |
| 2 | 6.463 | 740 | 744 | 750 | rBV | 128330 | 120279 | 5.26% | 0.389% |
| 3 | 6.610 | 766 | 769 | 772 | rBV | 148312 | 116120 | 5.08% | 0.376% |
| 4 | 6.845 | 805 | 809 | 812 | rBV | 825161 | 628887 | 27.50% | 2.036% |
| 5 | 6.998 | 832 | 835 | 839 | rBV | 100555 | 84638 | 3.70% | 0.274% |
| 6 | 7.398 | 898 | 903 | 908 | rBV | 87377 | 74644 | 3.26% | 0.242% |
| 7 | 8.128 | 1023 | 1027 | 1030 | rBV | 1101654 | 874366 | 38.23% | 2.831% |
| 8 | 9.198 | 1206 | 1209 | 1212 | rBV | 156778 | 114832 | 5.02% | 0.372% |
| 9 | 9.880 | 1321 | 1325 | 1328 | rBV | 1224442 | 980943 | 42.89% | 3.176% |
| 10 | 10.657 | 1454 | 1457 | 1462 | rBV3 | 60101 | 58181 | 2.54% | 0.188% |
| 11 | 11.233 | 1547 | 1555 | 1558 | rBV5 | 42956 | 57246 | 2.50% | 0.185% |
| 12 | 11.363 | 1573 | 1577 | 1579 | rBV | 1077452 | 1010633 | 44.19% | 3.272% |
| 13 | 11.380 | 1579 | 1580 | 1583 | rVV | 124578 | 89803 | 3.93% | 0.291% |
| 14 | 11.863 | 1658 | 1662 | 1664 | rVV | 144868 | 130500 | 5.71% | 0.423% |
| 15 | 11.886 | 1664 | 1666 | 1670 | rVB | 201910 | 171934 | 7.52% | 0.557% |
| 16 | 11.933 | 1671 | 1674 | 1677 | rBV | 89172 | 82596 | 3.61% | 0.267% |
| 17 | 11.968 | 1677 | 1680 | 1683 | rBV3 | 71099 | 70782 | 3.09% | 0.229% |
| 18 | 12.227 | 1721 | 1724 | 1728 | rVB | 53628 | 50931 | 2.23% | 0.165% |
| 19 | 12.304 | 1732 | 1737 | 1740 | rBV | 144839 | 141584 | 6.19% | 0.458% |
| 20 | 12.333 | 1740 | 1742 | 1744 | rBV | 114656 | 96694 | 4.23% | 0.313% |
| 21 | 12.357 | 1744 | 1746 | 1750 | rVB | 91133 | 76751 | 3.36% | 0.249% |
| 22 | 12.410 | 1751 | 1755 | 1758 | rBV | 130525 | 150563 | 6.58% | 0.487% |
| 23 | 12.451 | 1758 | 1762 | 1766 | rVB2 | 116605 | 166577 | 7.28% | 0.539% |
| 24 | 12.492 | 1766 | 1769 | 1773 | rBV | 100655 | 101764 | 4.45% | 0.329% |
| 25 | 12.568 | 1779 | 1782 | 1785 | rBV2 | 154266 | 147155 | 6.43% | 0.476% |
| 26 | 12.686 | 1800 | 1802 | 1806 | rVB2 | 74201 | 66996 | 2.93% | 0.217% |
| 27 | 12.727 | 1806 | 1809 | 1811 | rBV3 | 56559 | 60429 | 2.64% | 0.196% |
| 28 | 12.751 | 1811 | 1813 | 1816 | rVB | 71104 | 54651 | 2.39% | 0.177% |
| 29 | 12.798 | 1816 | 1821 | 1823 | rBV | 435490 | 417608 | 18.26% | 1.352% |
| 30 | 12.821 | 1823 | 1825 | 1828 | rVB2 | 131052 | 106582 | 4.66% | 0.345% |
| 31 | 12.857 | 1828 | 1831 | 1834 | rVB2 | 96173 | 106908 | 4.67% | 0.346% |
| 32 | 12.904 | 1834 | 1839 | 1842 | rBV4 | 59752 | 100941 | 4.41% | 0.327% |
| 33 | 12.933 | 1842 | 1844 | 1853 | rVB3 | 191600 | 267375 | 11.69% | 0.866% |
| 34 | 13.015 | 1854 | 1858 | 1865 | rBV3 | 137700 | 212387 | 9.29% | 0.688% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

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-BF083019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 13.068 | 1865 | 1867 | 1869 | rBV2 | 62350 | 56953 | 2.49% | 0.184% |
| 36 | 13.098 | 1869 | 1872 | 1874 | rBV3 | 65865 | 78363 | 3.43% | 0.254% |
| 37 | 13.121 | 1874 | 1876 | 1880 | rVB | 247167 | 239672 | 10.48% | 0.776% |
| 38 | 13.157 | 1880 | 1882 | 1884 | rBV2 | 47491 | 47980 | 2.10% | 0.155% |
| 39 | 13.192 | 1884 | 1888 | 1890 | rBV2 | 252279 | 250157 | 10.94% | 0.810% |
| 40 | 13.227 | 1890 | 1894 | 1897 | rBV | 507046 | 466613 | 20.40% | 1.511% |
| 41 | 13.257 | 1897 | 1899 | 1902 | rVV | 57557 | 70804 | 3.10% | 0.229% |
| 42 | 13.286 | 1902 | 1904 | 1908 | rVB2 | 61193 | 57655 | 2.52% | 0.187% |
| 43 | 13.321 | 1908 | 1910 | 1912 | rVB | 102033 | 68786 | 3.01% | 0.223% |
| 44 | 13.351 | 1912 | 1915 | 1920 | rBV4 | 102142 | 144531 | 6.32% | 0.468% |
| 45 | 13.427 | 1926 | 1928 | 1933 | rVB4 | 72782 | 115963 | 5.07% | 0.375% |
| 46 | 13.480 | 1933 | 1937 | 1939 | rBV3 | 83954 | 106695 | 4.66% | 0.345% |
| 47 | 13.527 | 1941 | 1945 | 1947 | rBV2 | 233271 | 292000 | 12.77% | 0.945% |
| 48 | 13.557 | 1947 | 1950 | 1955 | rVB | 315052 | 446226 | 19.51% | 1.445% |
| 49 | 13.610 | 1955 | 1959 | 1961 | rBV3 | 173887 | 254788 | 11.14% | 0.825% |
| 50 | 13.633 | 1961 | 1963 | 1964 | rBV | 153551 | 117152 | 5.12% | 0.379% |
| 51 | 13.698 | 1972 | 1974 | 1977 | rVB | 95077 | 74734 | 3.27% | 0.242% |
| 52 | 13.745 | 1977 | 1982 | 1985 | rBV4 | 156245 | 300595 | 13.14% | 0.973% |
| 53 | 13.815 | 1991 | 1994 | 1996 | rBV2 | 48966 | 54718 | 2.39% | 0.177% |
| 54 | 13.845 | 1996 | 1999 | 2002 | rVV3 | 123308 | 137715 | 6.02% | 0.446% |
| 55 | 13.933 | 2006 | 2014 | 2016 | rBV3 | 172272 | 351707 | 15.38% | 1.139% |
| 56 | 13.992 | 2016 | 2024 | 2027 | rVV2 | 1183502 | 2287144 | 100.00% | 7.405% |
| 57 | 14.021 | 2027 | 2029 | 2031 | rVV | 1354701 | 1084992 | 47.44% | 3.513% |
| 58 | 14.045 | 2031 | 2033 | 2037 | rVV | 184547 | 223287 | 9.76% | 0.723% |
| 59 | 14.080 | 2037 | 2039 | 2042 | rVB2 | 222305 | 174086 | 7.61% | 0.564% |
| 60 | 14.162 | 2051 | 2053 | 2059 | rVB4 | 163595 | 188455 | 8.24% | 0.610% |
| 61 | 14.210 | 2059 | 2061 | 2063 | rBV2 | 56523 | 49924 | 2.18% | 0.162% |
| 62 | 14.280 | 2069 | 2073 | 2076 | rBV4 | 105440 | 139453 | 6.10% | 0.452% |
| 63 | 14.345 | 2081 | 2084 | 2086 | rBV | 335490 | 393647 | 17.21% | 1.275% |
| 64 | 14.386 | 2086 | 2091 | 2093 | rVV | 1268442 | 1701641 | 74.40% | 5.510% |
| 65 | 14.415 | 2093 | 2096 | 2099 | rVV | 988021 | 897677 | 39.25% | 2.907% |
| 66 | 14.468 | 2099 | 2105 | 2108 | rVV5 | 393507 | 650761 | 28.45% | 2.107% |
| 67 | 14.504 | 2108 | 2111 | 2113 | rVV | 225320 | 233065 | 10.19% | 0.755% |
| 68 | 14.527 | 2113 | 2115 | 2118 | rVV | 251723 | 230763 | 10.09% | 0.747% |
| 69 | 14.574 | 2120 | 2123 | 2126 | rVV | 198278 | 201696 | 8.82% | 0.653% |
| 70 | 14.621 | 2129 | 2131 | 2134 | rBV3 | 65702 | 69153 | 3.02% | 0.224% |
| 71 | 14.668 | 2138 | 2139 | 2142 | rVV2 | 64408 | 55381 | 2.42% | 0.179% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

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-BF083019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 14.715 | 2143 | 2147 | 2153 | rVV | 349673 | 679067 | 29.69% | 2.199% |
| 73 | 14.768 | 2153 | 2156 | 2158 | rVV2 | 486154 | 516624 | 22.59% | 1.673% |
| 74 | 14.798 | 2158 | 2161 | 2165 | rVB2 | 358675 | 433312 | 18.95% | 1.403% |
| 75 | 14.874 | 2171 | 2174 | 2177 | rVB2 | 173754 | 176459 | 7.72% | 0.571% |
| 76 | 14.927 | 2179 | 2183 | 2189 | rVB3 | 336962 | 504600 | 22.06% | 1.634% |
| 77 | 15.027 | 2195 | 2200 | 2206 | rVV | 471002 | 623306 | 27.25% | 2.018% |
| 78 | 15.109 | 2210 | 2214 | 2220 | rVB3 | 186508 | 389996 | 17.05% | 1.263% |
| 79 | 15.327 | 2247 | 2251 | 2256 | rVB | 845163 | 1002750 | 43.84% | 3.247% |
| 80 | 15.386 | 2256 | 2261 | 2268 | rVB | 606813 | 798304 | 34.90% | 2.585% |
| 81 | 15.451 | 2268 | 2272 | 2275 | rBV | 654477 | 774665 | 33.87% | 2.508% |
| 82 | 15.486 | 2275 | 2278 | 2284 | rVB4 | 215399 | 403944 | 17.66% | 1.308% |
| 83 | 15.621 | 2297 | 2301 | 2304 | rBV3 | 136081 | 236220 | 10.33% | 0.765% |
| 84 | 15.768 | 2322 | 2326 | 2330 | rBV | 418673 | 607437 | 26.56% | 1.967% |
| 85 | 15.809 | 2330 | 2333 | 2341 | rVB | 499118 | 625583 | 27.35% | 2.026% |
| 86 | 15.886 | 2341 | 2346 | 2349 | rBV2 | 186082 | 312821 | 13.68% | 1.013% |
| 87 | 15.921 | 2349 | 2352 | 2356 | rVB2 | 163589 | 203695 | 8.91% | 0.660% |
| 88 | 15.968 | 2357 | 2360 | 2363 | rVB2 | 66108 | 85955 | 3.76% | 0.278% |
| 89 | 16.051 | 2370 | 2374 | 2377 | rVB2 | 93417 | 110699 | 4.84% | 0.358% |
| 90 | 16.151 | 2388 | 2391 | 2394 | rBV2 | 69129 | 76652 | 3.35% | 0.248% |
| 91 | 16.256 | 2405 | 2409 | 2413 | rBV2 | 165663 | 240956 | 10.54% | 0.780% |
| 92 | 16.709 | 2482 | 2486 | 2490 | rVB2 | 91199 | 139784 | 6.11% | 0.453% |
| 93 | 16.898 | 2513 | 2518 | 2525 | rVB2 | 279939 | 618506 | 27.04% | 2.003% |
| 94 | 17.074 | 2544 | 2548 | 2552 | rVB | 101371 | 158239 | 6.92% | 0.512% |
| 95 | 17.127 | 2552 | 2557 | 2563 | rBV | 114475 | 194484 | 8.50% | 0.630% |
| 96 | 17.333 | 2587 | 2592 | 2601 | rVB | 299861 | 640263 | 27.99% | 2.073% |
| 97 | 17.703 | 2648 | 2655 | 2657 | rBV3 | 102093 | 220272 | 9.63% | 0.713% |
| 98 | 17.821 | 2671 | 2675 | 2680 | rBV5 | 46214 | 87154 | 3.81% | 0.282% |
| 99 | 17.986 | 2699 | 2703 | 2714 | rVB5 | 72650 | 164280 | 7.18% | 0.532% |
| 100 | 18.098 | 2716 | 2722 | 2732 | rVB | 175407 | 444681 | 19.44% | 1.440% |

Sum of corrected areas: 30884778

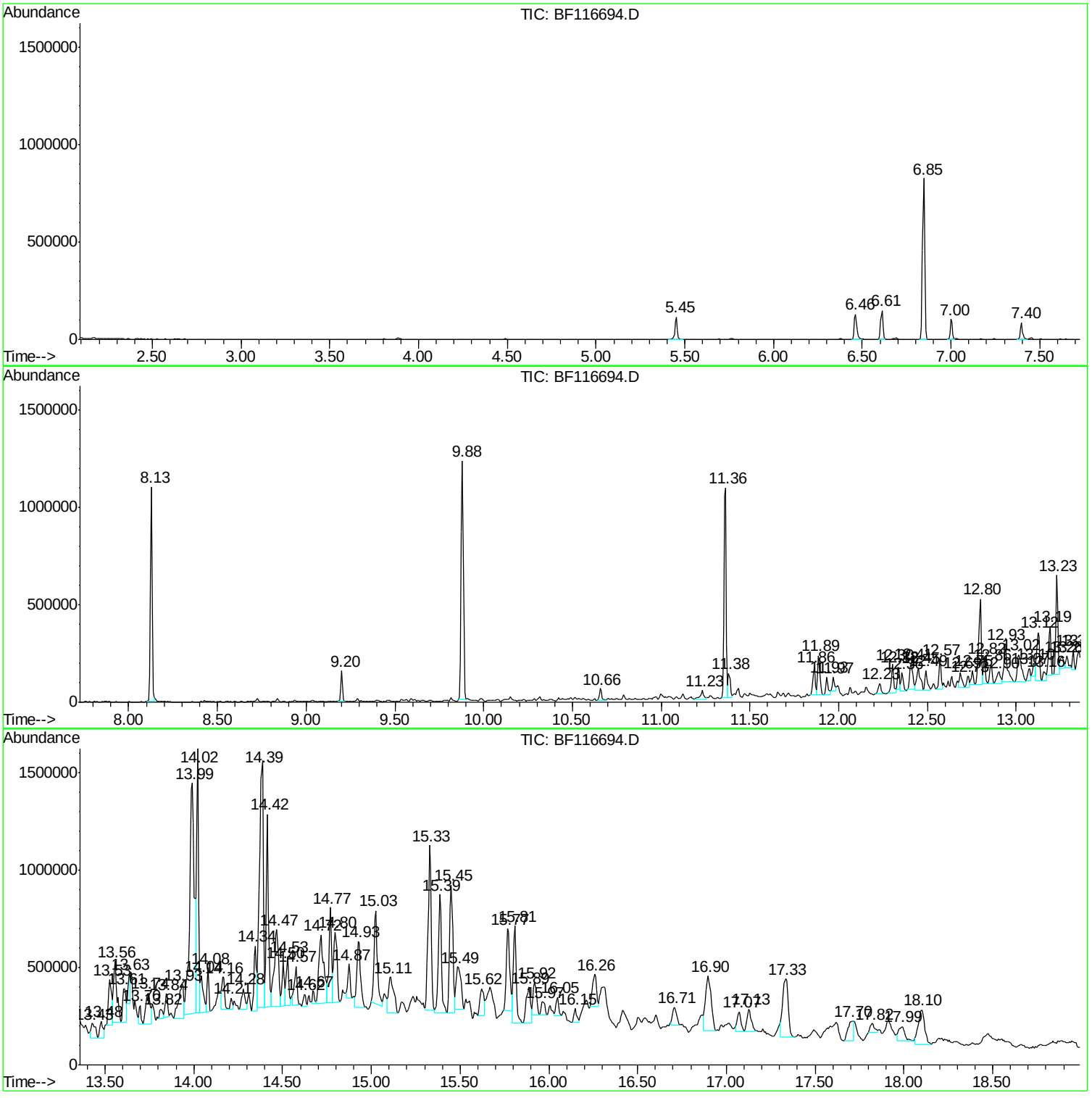
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

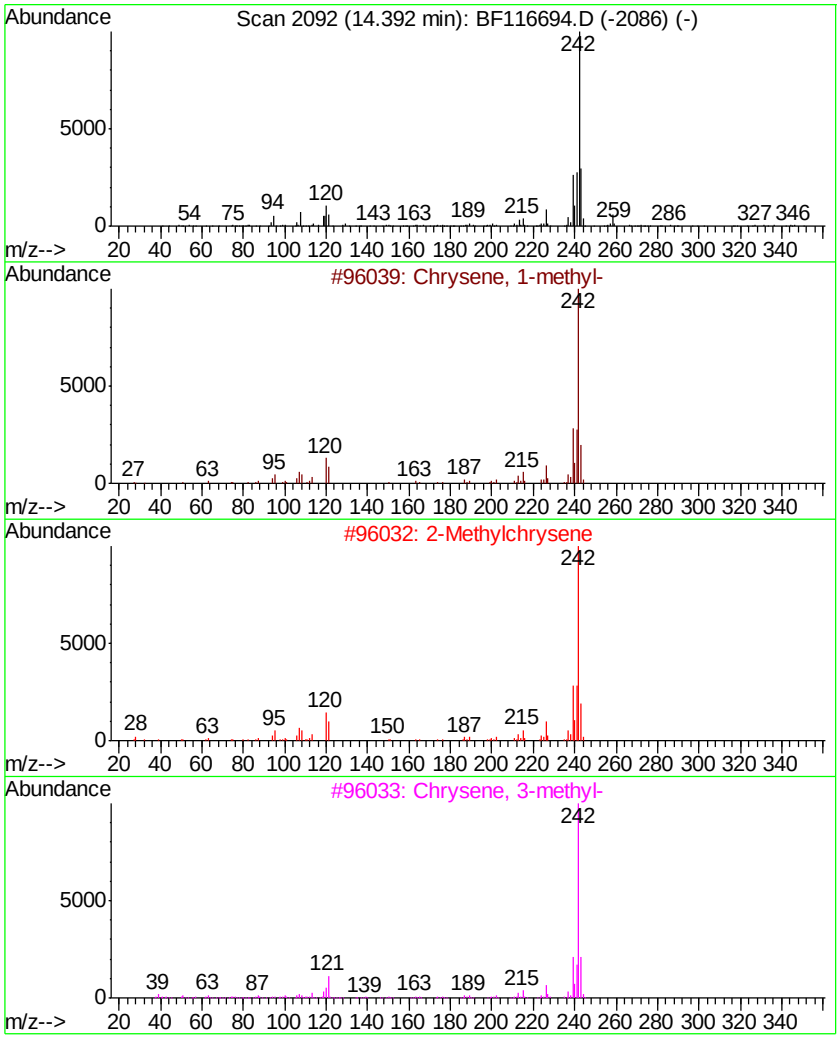
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

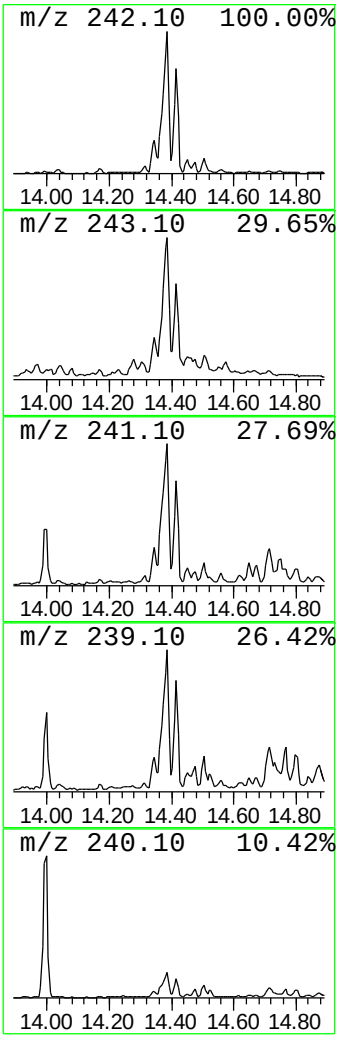
**********

Peak Number 1 Chrysene, 1-methyl- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.39 | 14.88 ng | 1701640 | Chrysene-d12 | 14.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Chrysene, 1-methyl- | 242 | C19H14 | 003351-28-8 | 96 |
| 2 | 2-Methylchrysene | 242 | C19H14 | 003351-32-4 | 96 |
| 3 | Chrysene, 3-methyl- | 242 | C19H14 | 003351-31-3 | 95 |
| 4 | Benz[a]anthracene, 7-methyl- | 242 | C19H14 | 002541-69-7 | 95 |
| 5 | Triphenylene, 2-methyl- | 242 | C19H14 | 001705-84-6 | 95 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

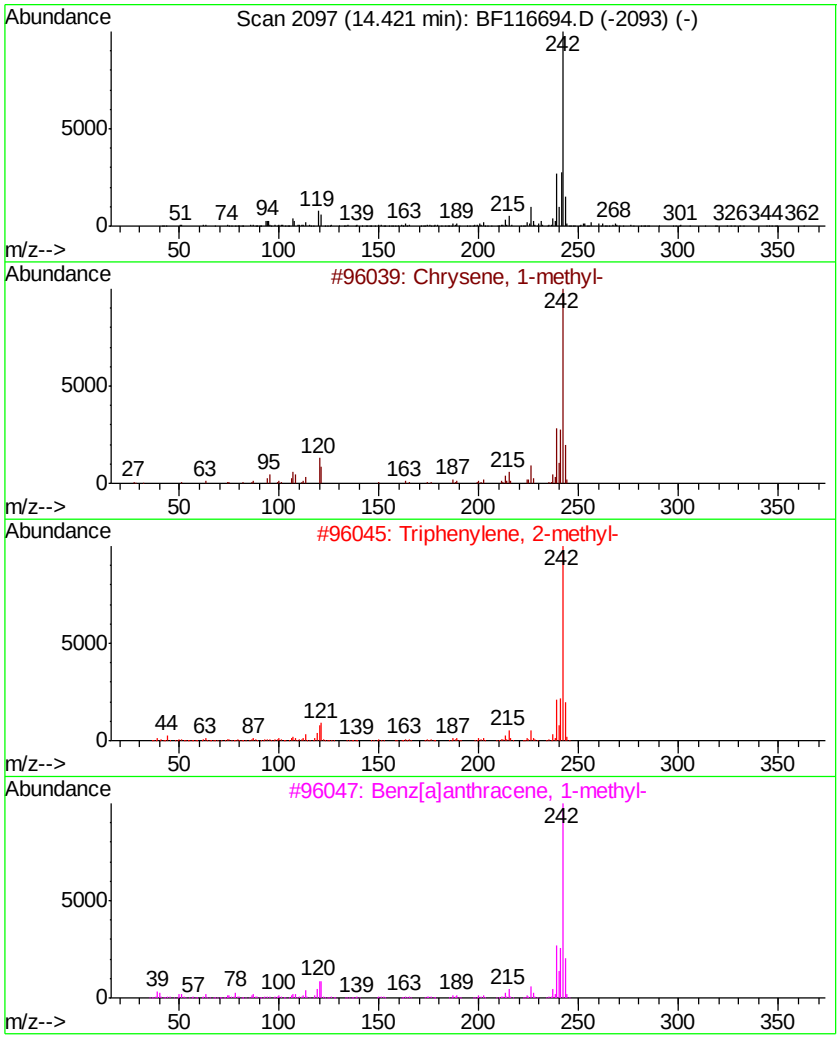
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

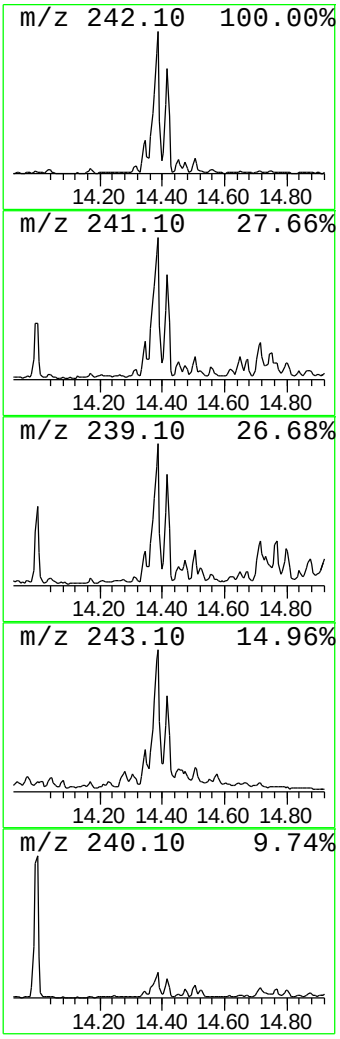
**********

Peak Number 2 Triphenylene, 2-methyl- Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.42 | 7.85 ng | 897677 | Chrysene-d12 | 14.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Chrysene, 1-methyl- | 242 | C19H14 | 003351-28-8 | 96 |
| 2 | Triphenylene, 2-methyl- | 242 | C19H14 | 001705-84-6 | 93 |
| 3 | Benz[a]anthracene, 1-methyl- | 242 | C19H14 | 002498-77-3 | 93 |
| 4 | Benz[a]anthracene, 7-methyl- | 242 | C19H14 | 002541-69-7 | 90 |
| 5 | Chrysene, 5-methyl- | 242 | C19H14 | 003697-24-3 | 81 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

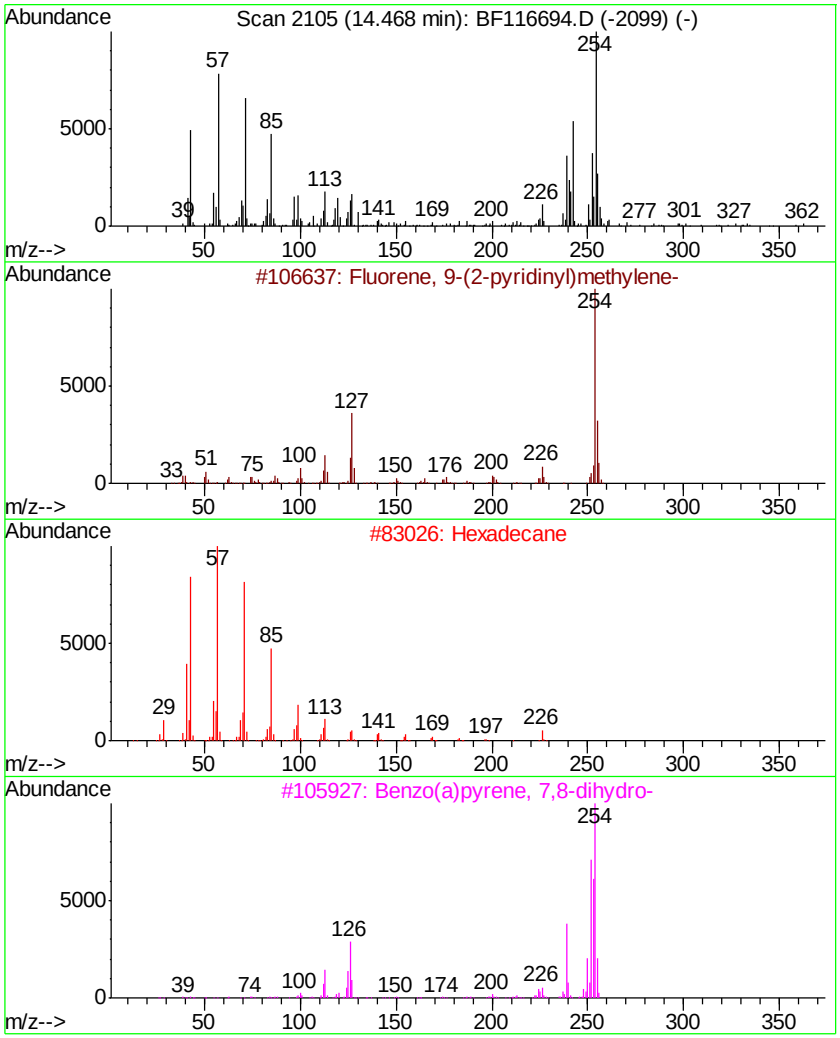
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

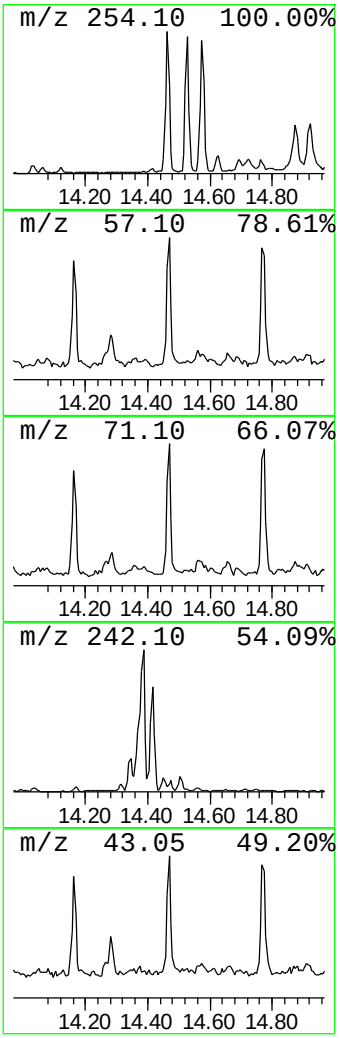
**********

Peak Number 3 unknown14.47 Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.47 | 5.69 ng | 650761 | Chrysene-d12 | 14.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fluorene, 9-(2-pyridinyl)methylene- | 255 | C19H13N | 002871-27-4 | 49 |
| 2 | Hexadecane | 226 | C16H34 | 000544-76-3 | 45 |
| 3 | Benzo(a)pyrene, 7,8-dihydro- | 254 | C20H14 | 017573-23-8 | 38 |
| 4 | Benzo(e)pyrene, 9,10-dihydro- | 254 | C20H14 | 066788-01-0 | 30 |
| 5 | 2,2'-Binaphthalene | 254 | C20H14 | 000612-78-2 | 30 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

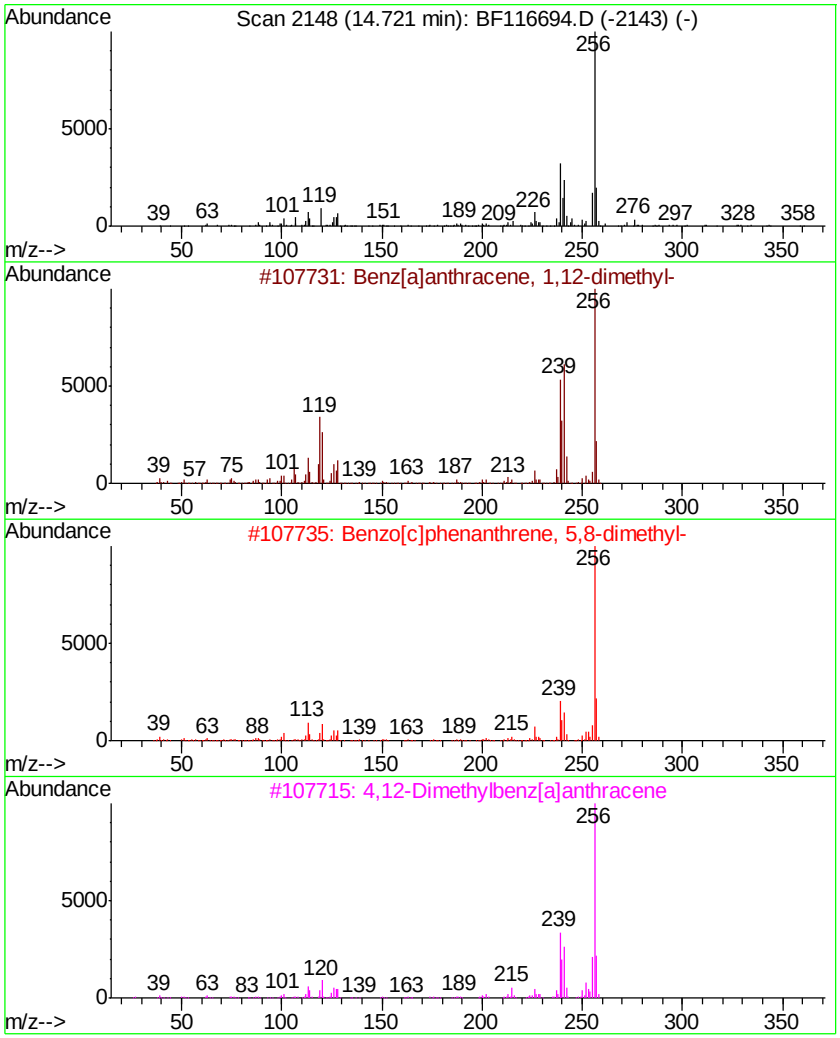
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

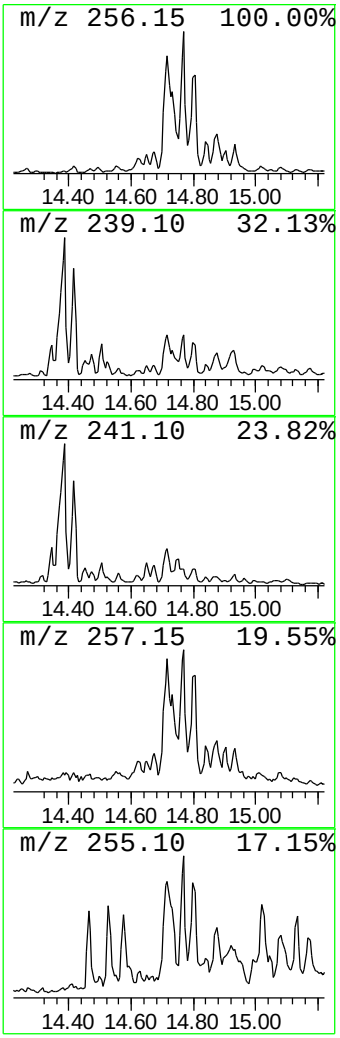
**********

Peak Number 4 Benz[a]anthracene, 1,12-dim... Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.72 | 5.94 ng | 679067 | Chrysene-d12 | 14.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benz[a]anthracene, 1,12-dimethyl- | 256 | C20H16 | 000313-74-6 | 95 |
| 2 | Benzo[c]phenanthrene, 5,8-dimethyl- | 256 | C20H16 | 054986-63-9 | 93 |
| 3 | 4,12-Dimethylbenz[a]anthracene | 256 | C20H16 | 035187-19-0 | 93 |
| 4 | Benz[a]anthracene, 7,12-dimethyl- | 256 | C20H16 | 000057-97-6 | 93 |
| 5 | 5,12-Dimethylbenz[a]anthracene | 256 | C20H16 | 021297-22-3 | 91 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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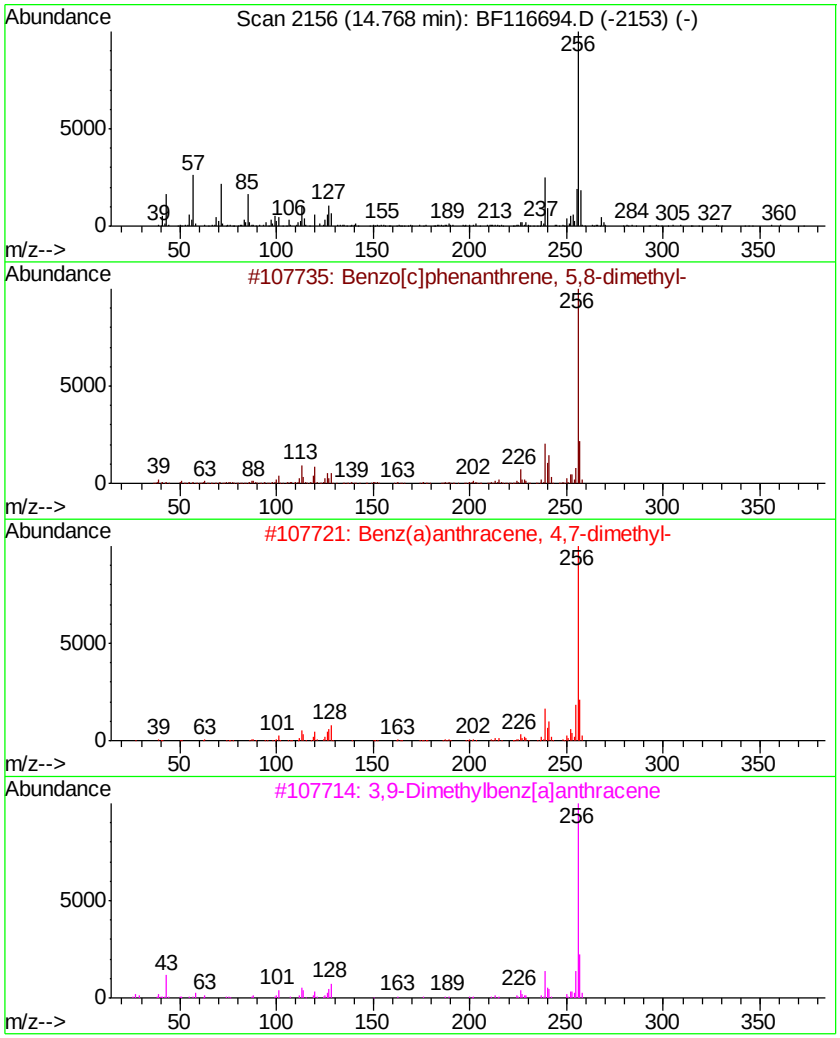
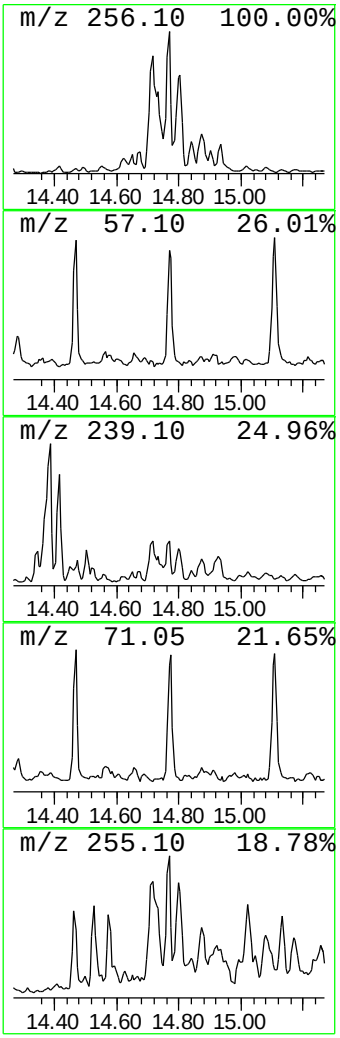
Peak Number 5 Benzo[c]phenanthrene, 5,8-d... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.77 | 13.34 ng | 516624 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[c]phenanthrene, 5,8-dimethyl- | 256 | C20H16 | 054986-63-9 | 93 |
| 2 | Benz(a)anthracene, 4,7-dimethyl- | 256 | C20H16 | 035187-18-9 | 93 |
| 3 | 3,9-Dimethylbenz[a]anthracene | 256 | C20H16 | 000316-51-8 | 91 |
| 4 | 5,12-Dimethylbenz[a]anthracene | 256 | C20H16 | 021297-22-3 | 81 |
| 5 | Benz(a)anthracene, 8,12-dimethyl- | 256 | C20H16 | 020627-31-0 | 81 |



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Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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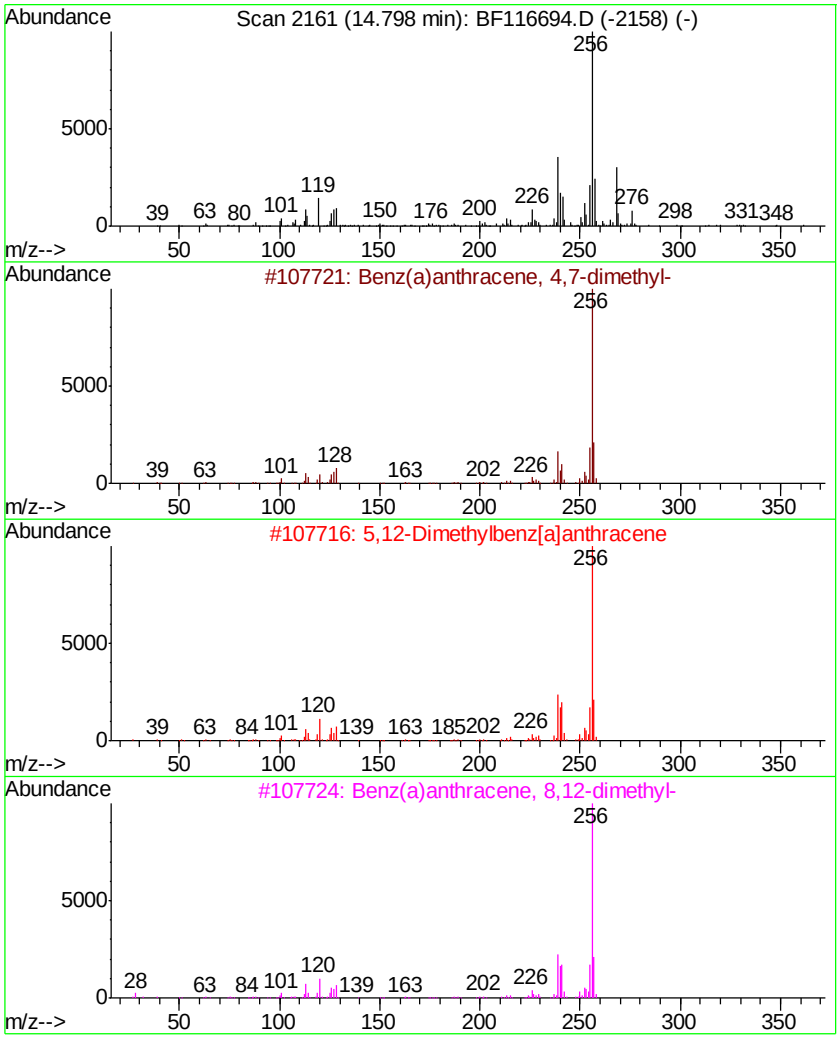
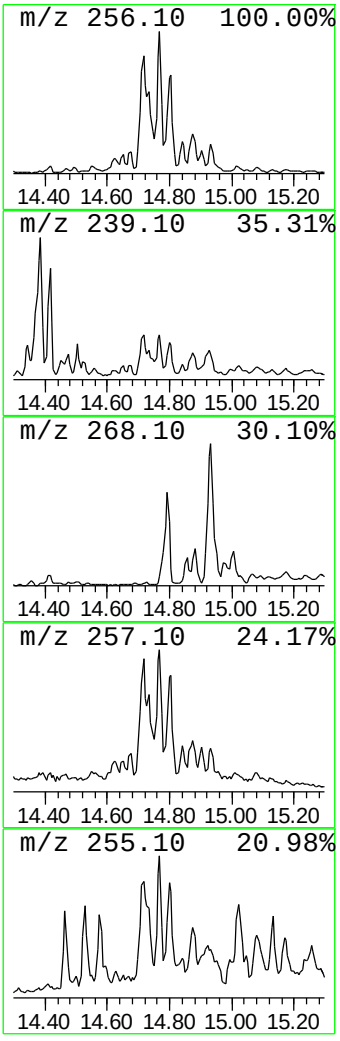
Peak Number 6 Benz(a)anthracene, 4,7-dime... Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.80 | 11.19 ng | 433312 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benz(a)anthracene, 4,7-dimethyl- | 256 | C20H16 | 035187-18-9 | 93 |
| 2 | 5,12-Dimethylbenz[a]anthracene | 256 | C20H16 | 021297-22-3 | 93 |
| 3 | Benz(a)anthracene, 8,12-dimethyl- | 256 | C20H16 | 020627-31-0 | 93 |
| 4 | Benz[a]anthracene, 7,12-dimethyl- | 256 | C20H16 | 000057-97-6 | 91 |
| 5 | 4,12-Dimethylbenz[a]anthracene | 256 | C20H16 | 035187-19-0 | 89 |



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Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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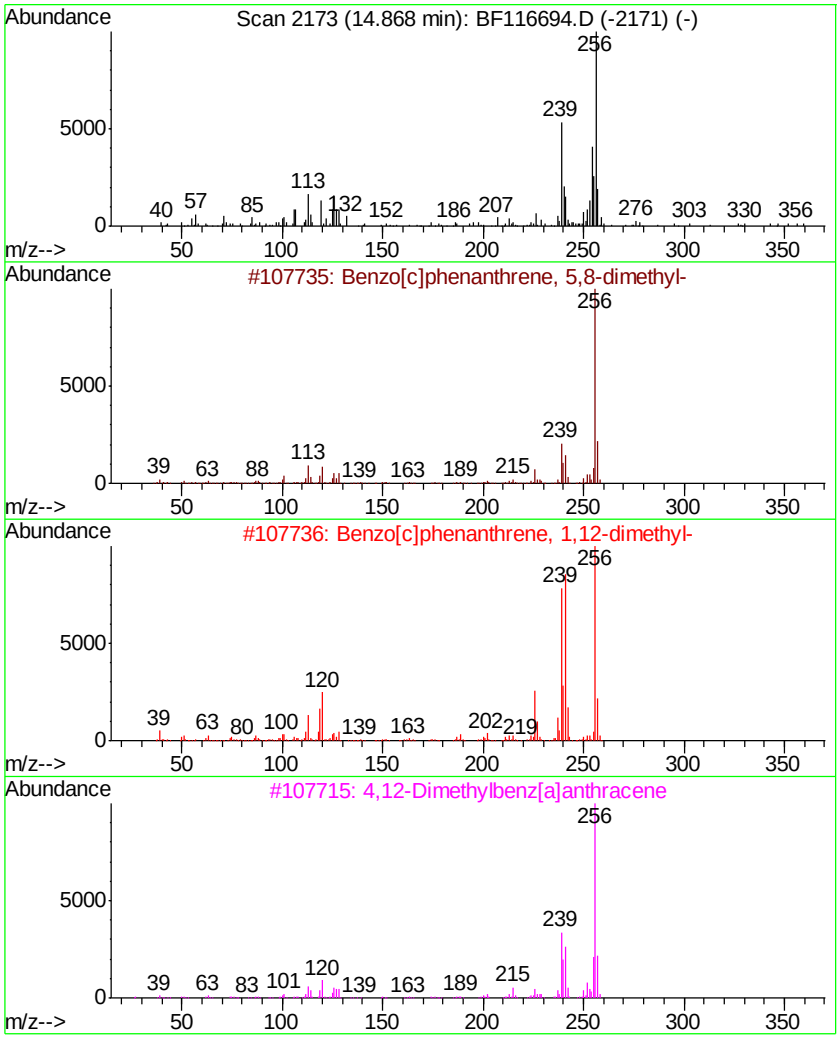
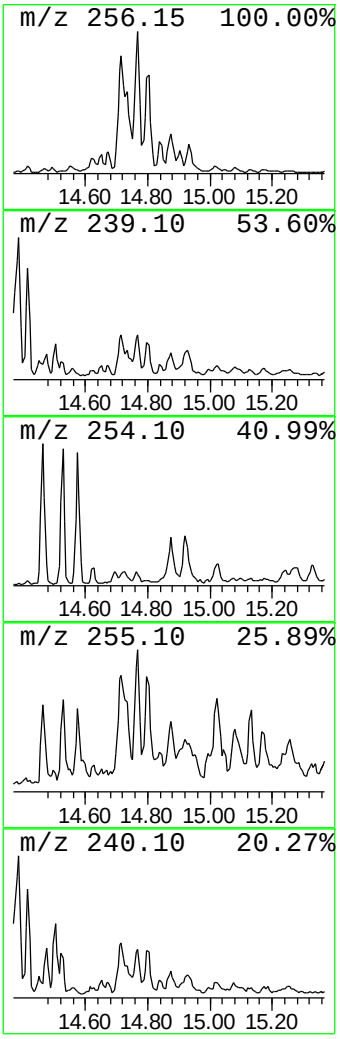
Peak Number 7 Benzo[c]phenanthrene, 1,12-... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.87 | 4.56 ng | 176459 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[c]phenanthrene, 5,8-dimethyl- | 256 | C20H16 | 054986-63-9 | 86 |
| 2 | Benzo[c]phenanthrene, 1,12-dimet... | 256 | C20H16 | 004076-43-1 | 83 |
| 3 | 4,12-Dimethylbenz[a]anthracene | 256 | C20H16 | 035187-19-0 | 78 |
| 4 | Benz[a]anthracene, 7,12-dimethyl- | 256 | C20H16 | 000057-97-6 | 78 |
| 5 | Benz(a)anthracene, 8,12-dimethyl- | 256 | C20H16 | 020627-31-0 | 70 |



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Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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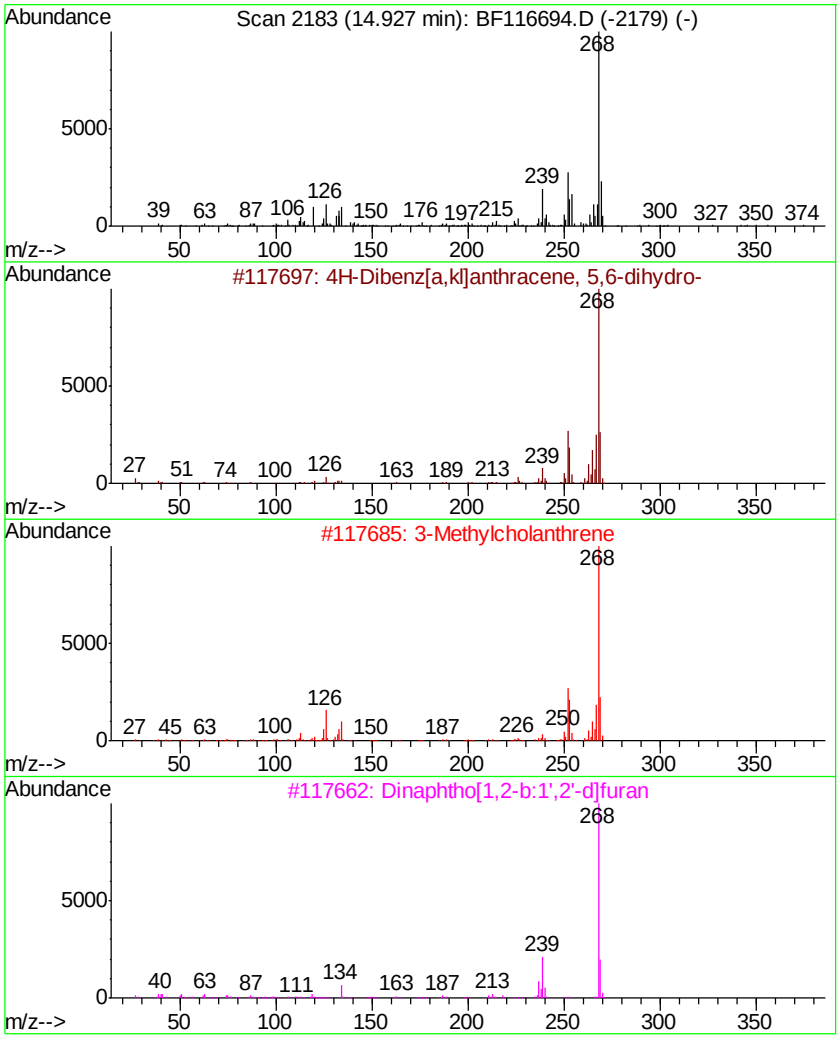
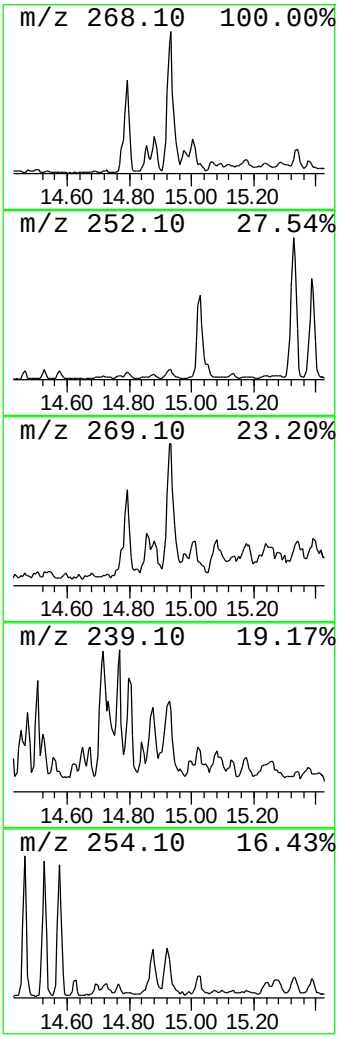
Peak Number 8 4H-Dibenz[a,kl]anthracene, ... Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.93 | 13.03 ng | 504600 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4H-Dibenz[a,kl]anthracene, 5,6-d... | 268 | C21H16 | 007198-87-0 | 76 |
| 2 | 3-Methylcholanthrene | 268 | C21H16 | 000056-49-5 | 62 |
| 3 | Dinaphtho[1,2-b:1',2'-d]furan | 268 | C20H12O | 000207-93-2 | 60 |
| 4 | 9,10-Dihydro-8-methylbenzo[a]pyrene | 268 | C21H16 | 089524-72-1 | 58 |
| 5 | Naphthalene, 1,1'-methylenebis- | 268 | C21H16 | 000607-50-1 | 58 |



# Library Search Compound Report

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Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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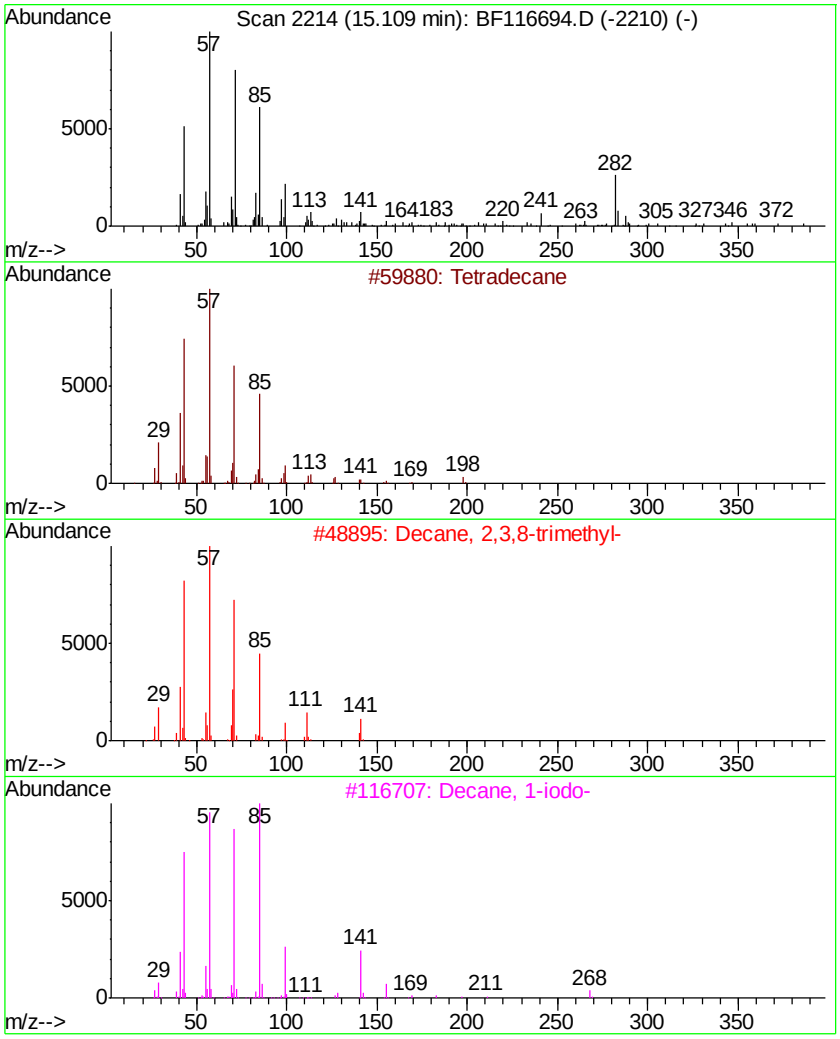
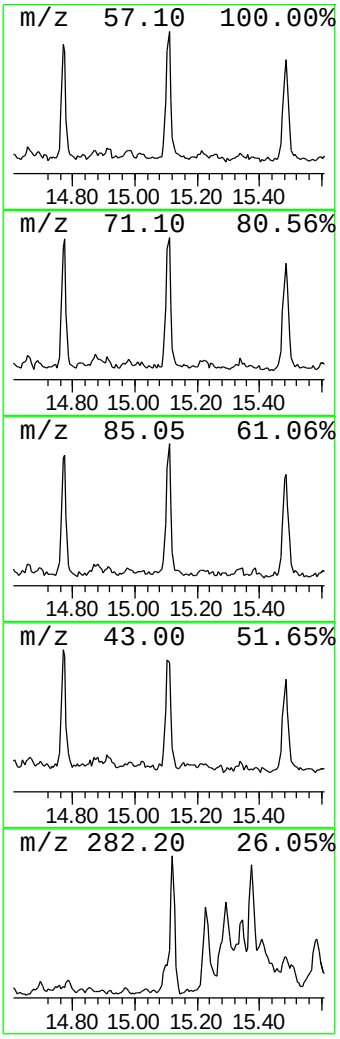
Peak Number 9 Tetradecane Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.11 | 10.07 ng | 389996 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tetradecane | 198 | C14H30 | 000629-59-4 | 59 |
| 2 | Decane, 2,3,8-trimethyl- | 184 | C13H28 | 062238-14-6 | 59 |
| 3 | Decane, 1-iodo- | 268 | C10H21I | 002050-77-3 | 53 |
| 4 | Undecane, 2,9-dimethyl- | 184 | C13H28 | 017301-26-7 | 53 |
| 5 | Decane, 5-propyl- | 184 | C13H28 | 017312-62-8 | 52 |



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Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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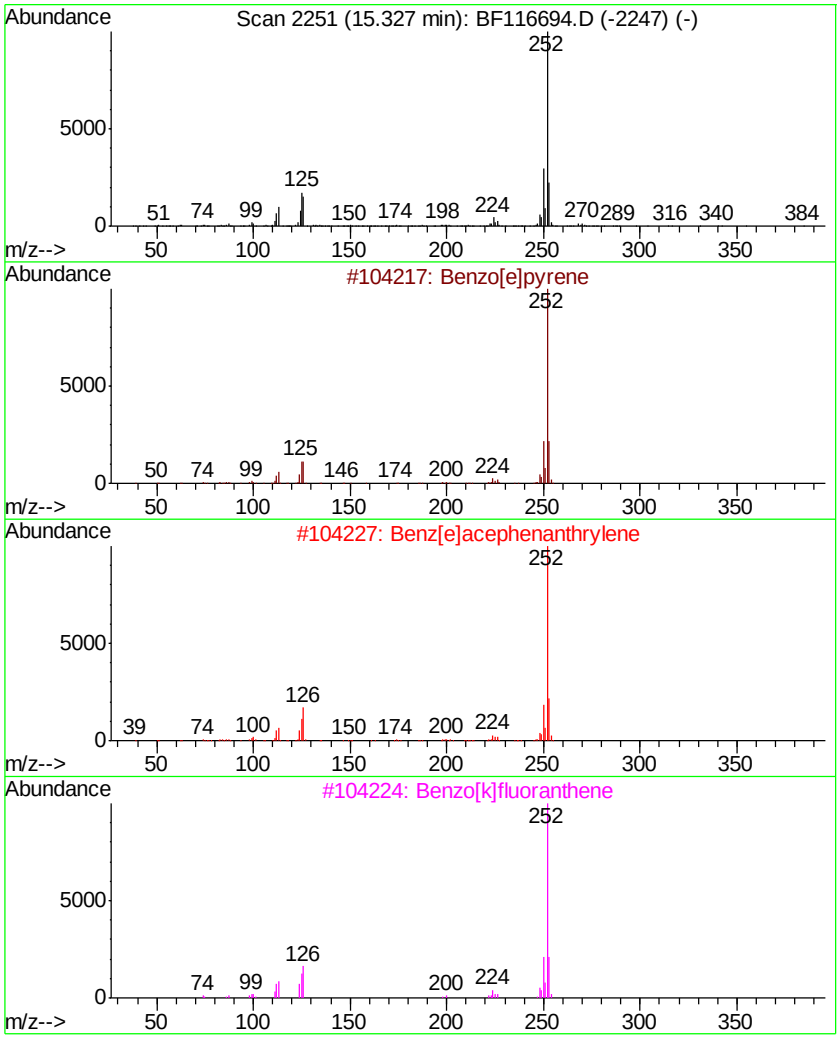
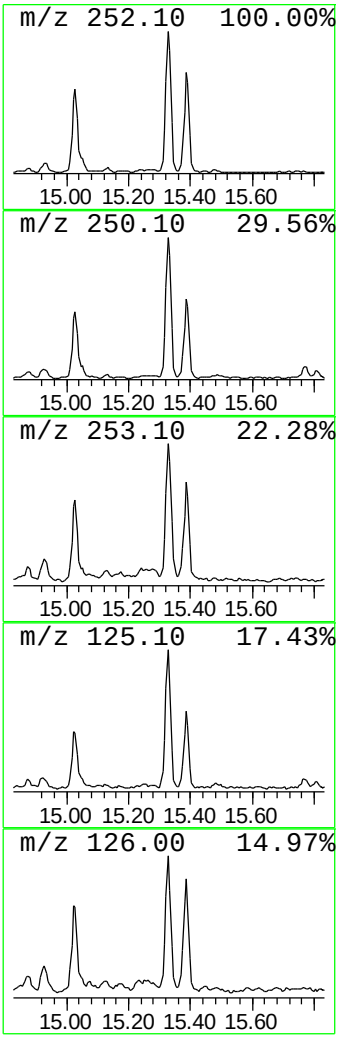
Peak Number 10 Benzo[e]pyrene Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.33 | 25.89 ng | 1002750 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 98 |
| 2 | Benz[e]acephenanthrylene | 252 | C20H12 | 000205-99-2 | 98 |
| 3 | Benzo[k]fluoranthene | 252 | C20H12 | 000207-08-9 | 98 |
| 4 | Perylene | 252 | C20H12 | 000198-55-0 | 96 |
| 5 | Benzo[a]pyrene | 252 | C20H12 | 000050-32-8 | 94 |



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Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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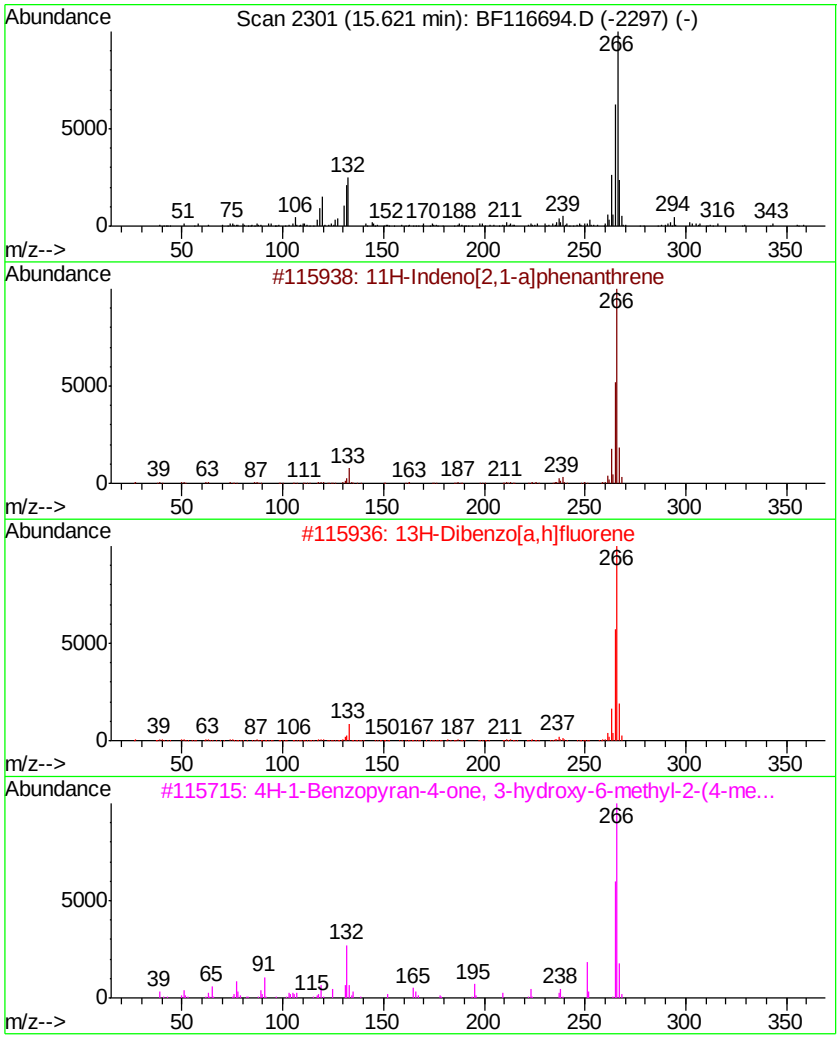
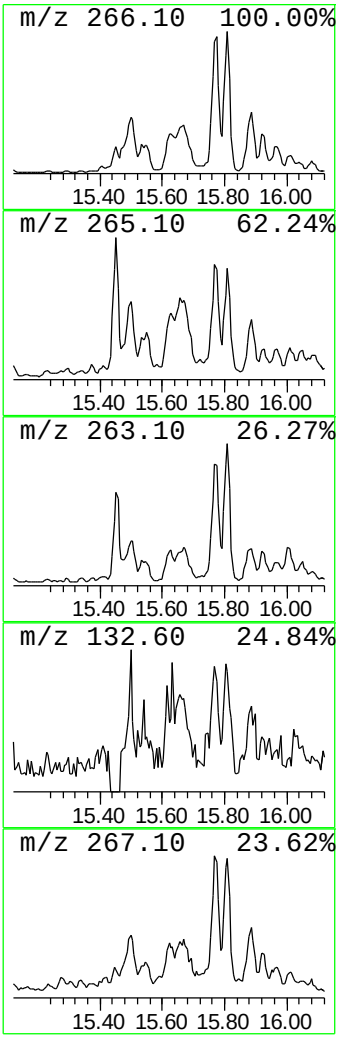
Peak Number 11 11H-Indeno[2,1-a]phenanthrene Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.62 | 6.10 ng | 236220 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 11H-Indeno[2,1-a]phenanthrene | 266 | C21H14 | 000220-97-3 | 64 |
| 2 | 13H-Dibenzo[a,h]fluorene | 266 | C21H14 | 000239-85-0 | 64 |
| 3 | 4H-1-Benzopyran-4-one, 3-hydroxy... | 266 | C17H14O3 | 078396-37-9 | 58 |
| 4 | Thiazole, 4-phenyl-2-(4-tolylami... | 266 | C16H14N2S | 016098-04-7 | 50 |
| 5 | 8H-Indeno[2,1-b]phenanthrene | 266 | C21H14 | 000241-28-1 | 49 |



Library Search Compound Report

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Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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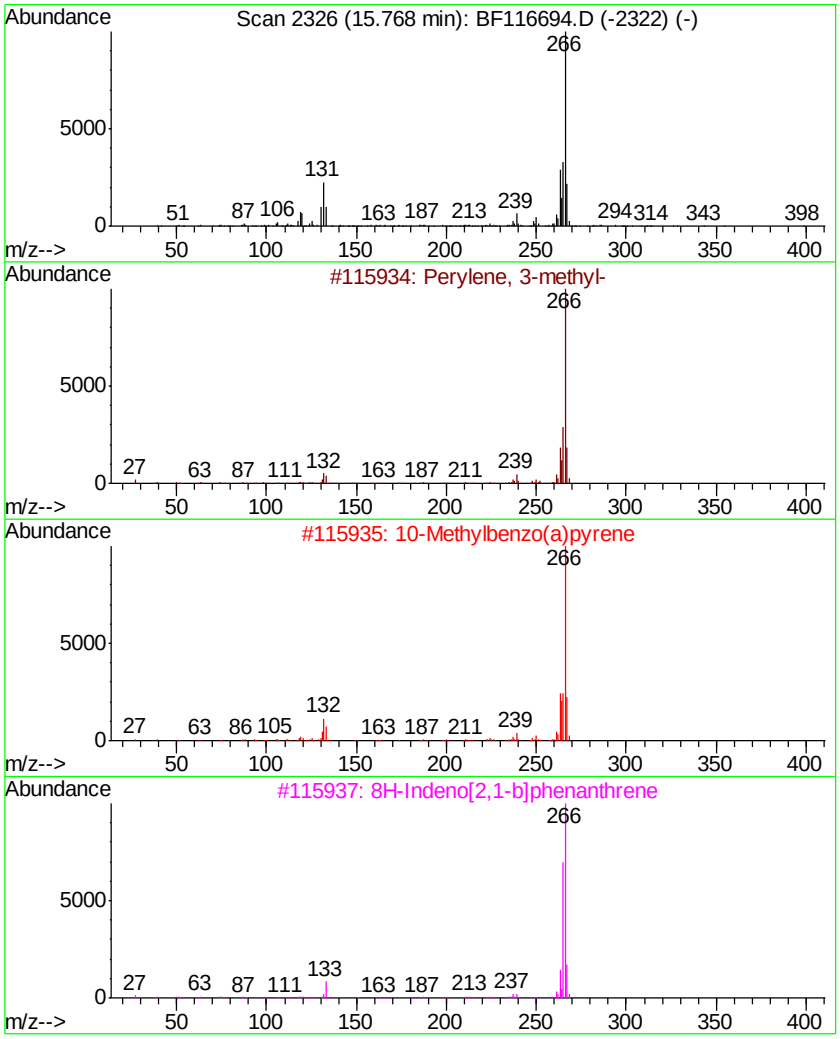
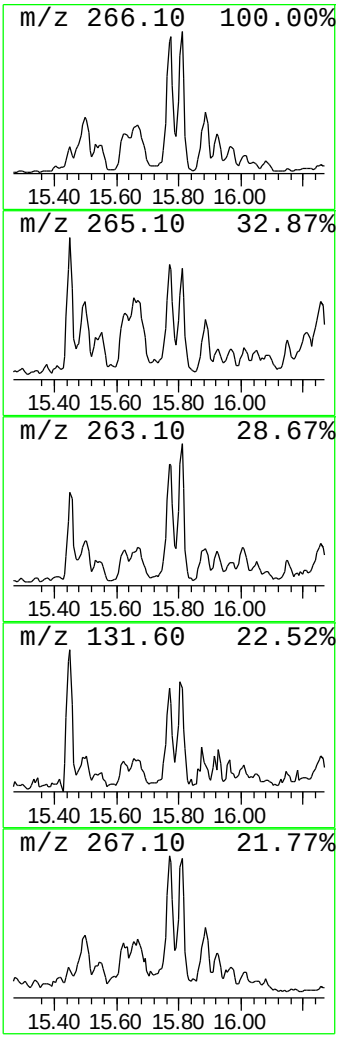
Peak Number 12 Perylene, 3-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.77 | 15.68 ng | 607437 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Perylene, 3-methyl- | 266 | C21H14 | 024471-47-4 | 94 |
| 2 | 10-Methylbenzo(a)pyrene | 266 | C21H14 | 063104-32-5 | 93 |
| 3 | 8H-Indeno[2,1-b]phenanthrene | 266 | C21H14 | 000241-28-1 | 70 |
| 4 | 13H-Dibenzo[a,h]fluorene | 266 | C21H14 | 000239-85-0 | 64 |
| 5 | Benz[j]aceanthrylene, 3-methyl- | 266 | C21H14 | 003343-10-0 | 64 |



Library Search Compound Report

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Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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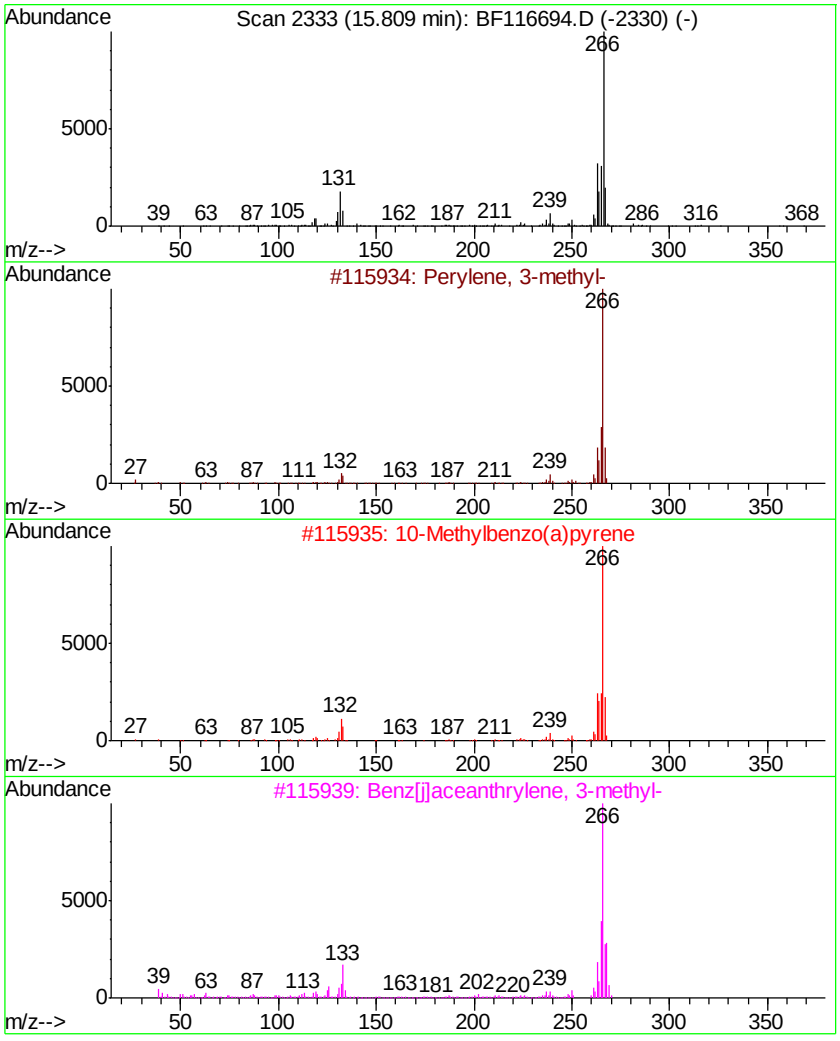
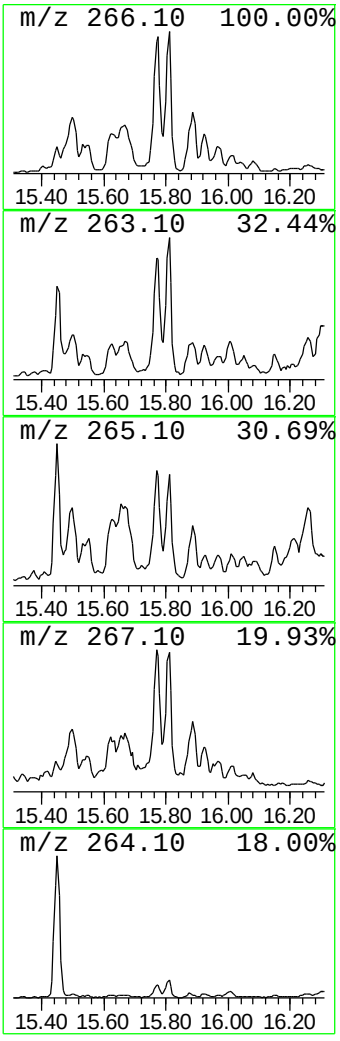
Peak Number 13 10-Methylbenzo(a)pyrene Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.81 | 16.15 ng | 625583 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Perylene, 3-methyl- | 266 | C21H14 | 024471-47-4 | 89 |
| 2 | 10-Methylbenzo(a)pyrene | 266 | C21H14 | 063104-32-5 | 86 |
| 3 | Benz[j]aceanthrylene, 3-methyl- | 266 | C21H14 | 003343-10-0 | 60 |
| 4 | 11H-Indeno[2,1-a]phenanthrene | 266 | C21H14 | 000220-97-3 | 53 |
| 5 | 8H-Indeno[2,1-b]phenanthrene | 266 | C21H14 | 000241-28-1 | 52 |



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Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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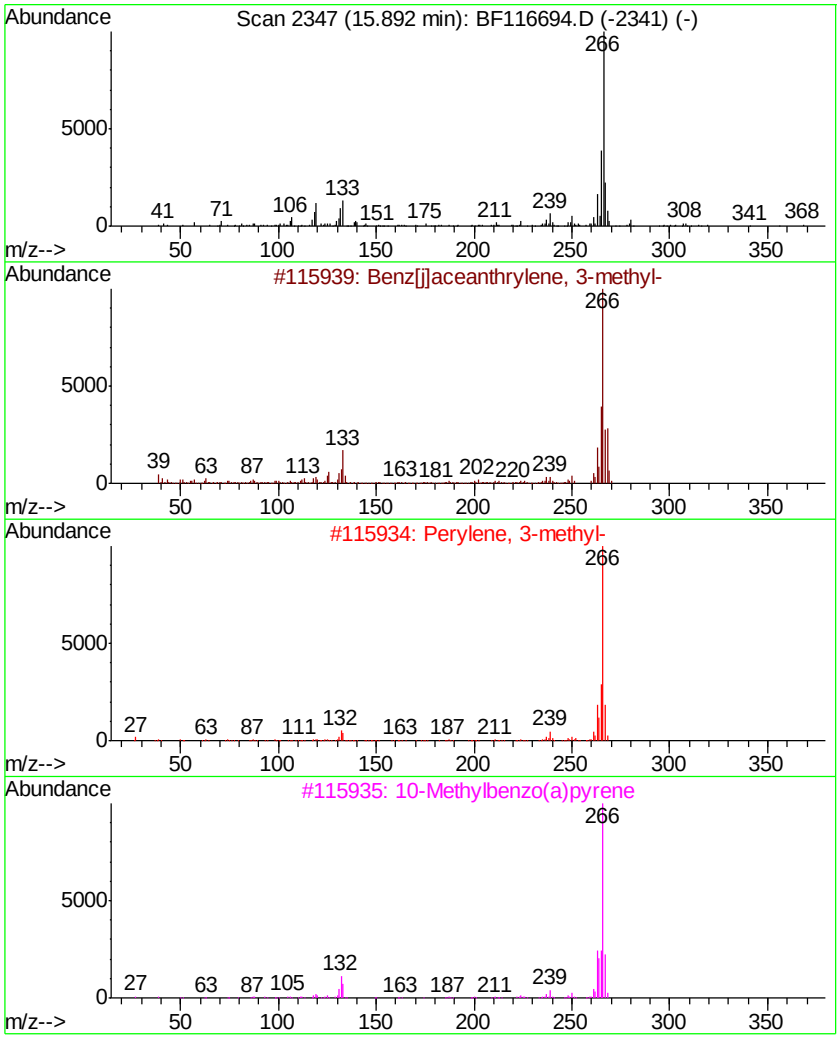
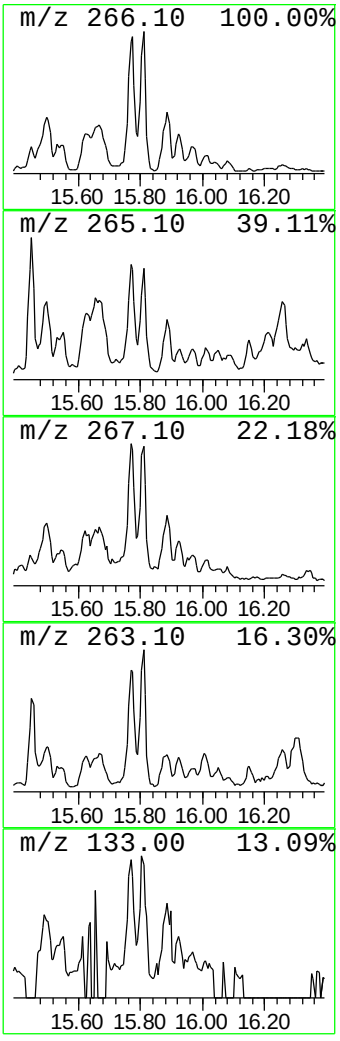
Peak Number 14 Benz[j]aceanthrylene, 3-met... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.89 | 8.08 ng | 312821 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benz[j]aceanthrylene, 3-methyl- | 266 | C21H14 | 003343-10-0 | 91 |
| 2 | Perylene, 3-methyl- | 266 | C21H14 | 024471-47-4 | 87 |
| 3 | 10-Methylbenzo(a)pyrene | 266 | C21H14 | 063104-32-5 | 76 |
| 4 | 8H-Indeno[2,1-b]phenanthrene | 266 | C21H14 | 000241-28-1 | 76 |
| 5 | 11H-Indeno[2,1-a]phenanthrene | 266 | C21H14 | 000220-97-3 | 70 |



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Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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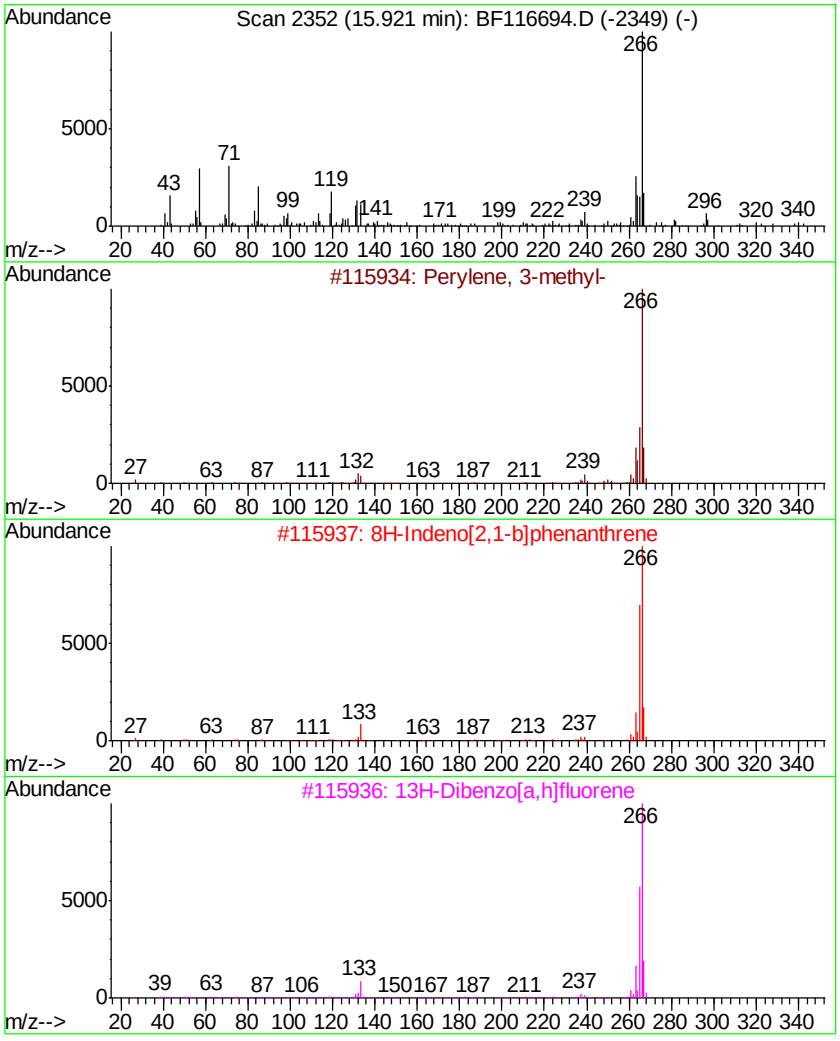
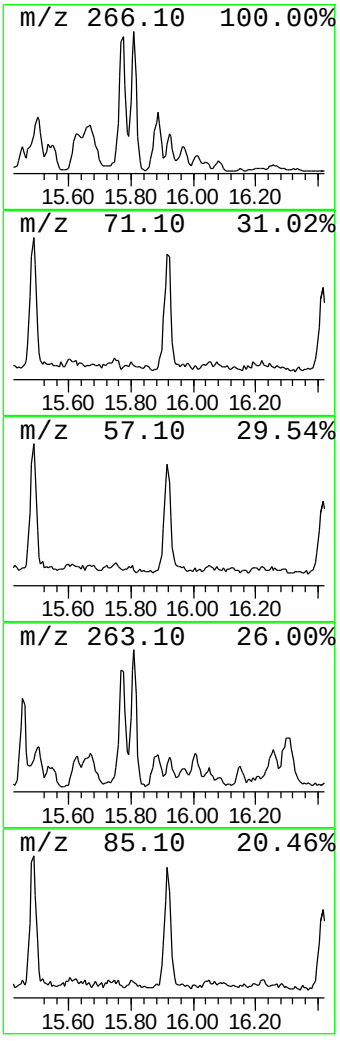
Peak Number 15 8H-Indeno[2,1-b]phenanthrene Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.92 | 5.26 ng | 203695 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Perylene, 3-methyl- | 266 | C21H14 | 024471-47-4 | 60 |
| 2 | 8H-Indeno[2,1-b]phenanthrene | 266 | C21H14 | 000241-28-1 | 58 |
| 3 | 13H-Dibenzo[a,h]fluorene | 266 | C21H14 | 000239-85-0 | 52 |
| 4 | 11H-Indeno[2,1-a]phenanthrene | 266 | C21H14 | 000220-97-3 | 52 |
| 5 | 9-Trifluoromethyl-5,6,7,8-tetrah... | 266 | C12H9F3N4 | 1000301-28-5 | 49 |



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Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

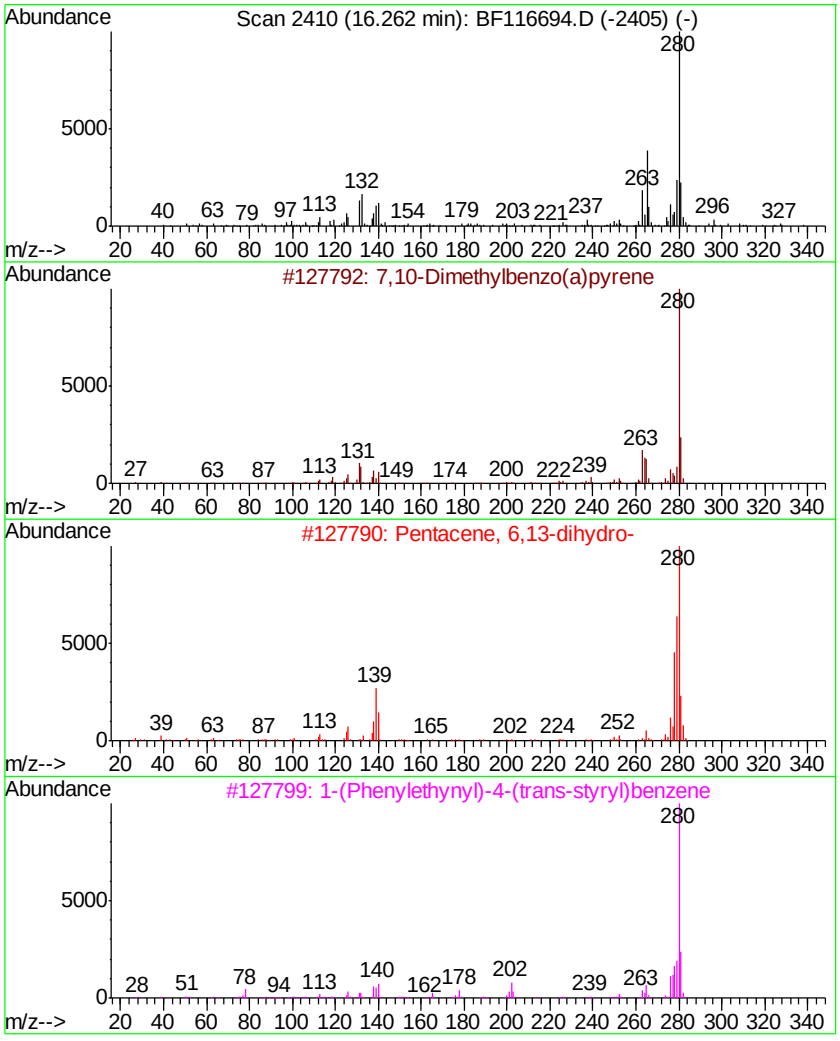
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

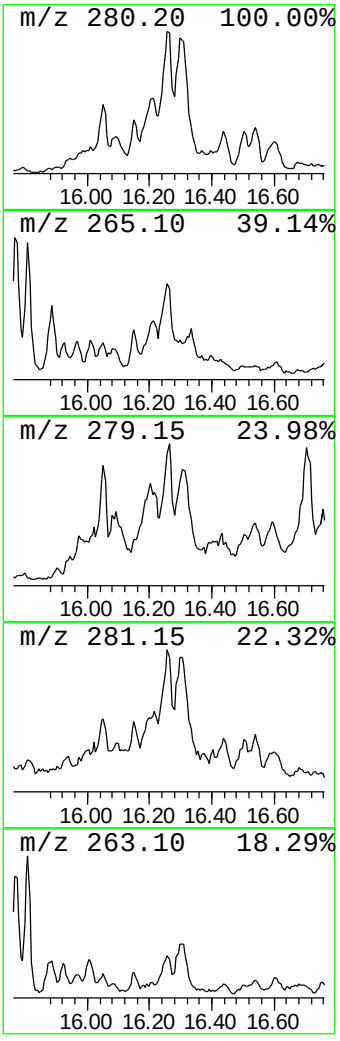
**********

Peak Number 16 7,10-Dimethylbenzo(a)pyrene Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.26 | 6.22 ng | 240956 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 7,10-Dimethylbenzo(a)pyrene | 280 | C22H16 | 063104-33-6 | 92 |
| 2 | Pentacene, 6,13-dihydro- | 280 | C22H16 | 013579-08-3 | 64 |
| 3 | 1-(Phenylethynyl)-4-(trans-styryl... | 280 | C22H16 | 173035-17-1 | 60 |
| 4 | Naphthalene, 1,7-diphenyl- | 280 | C22H16 | 000970-06-9 | 60 |
| 5 | Benzo[h]naphtho[1,2-c]cinnoline | 280 | C20H12N2 | 000213-51-4 | 60 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

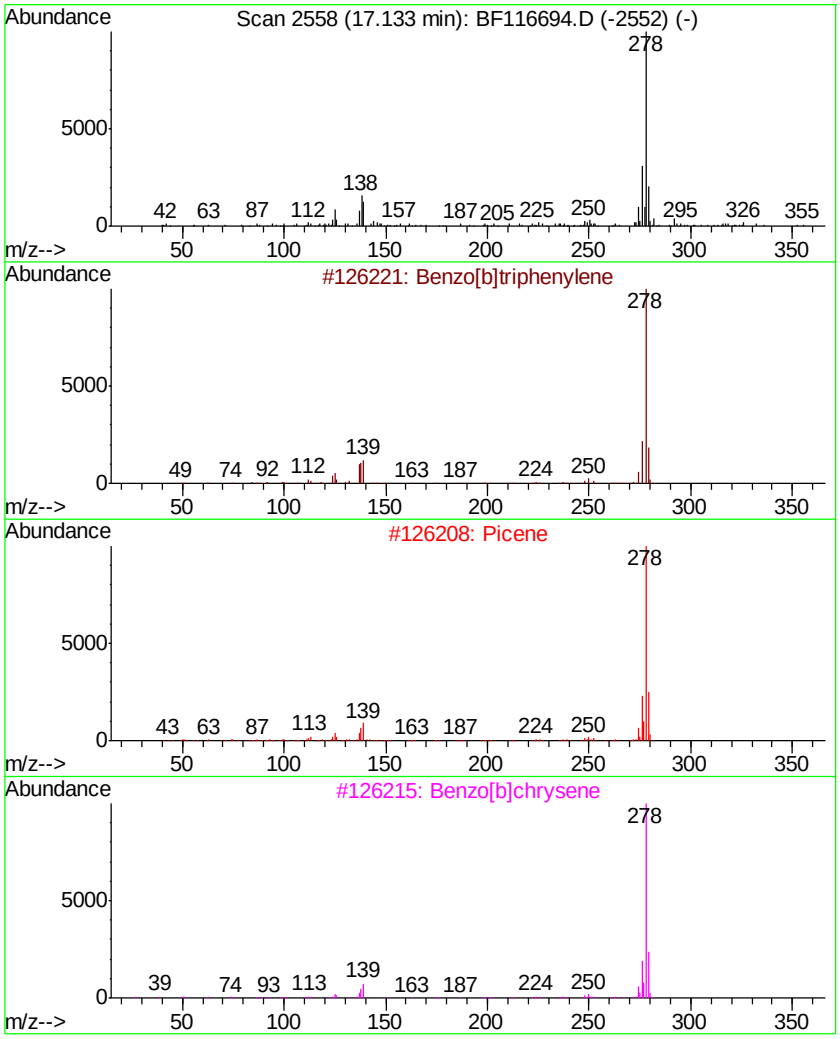
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

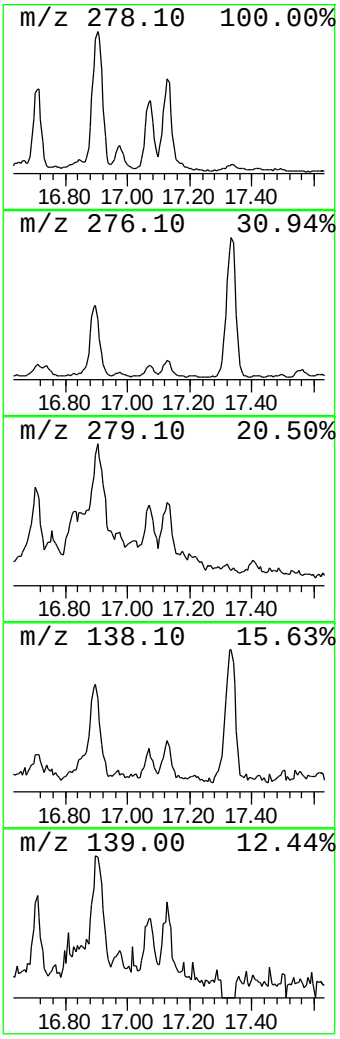
**********

Peak Number 17 Benzo[b]triphenylene Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.13 | 5.02 ng | 194484 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[b]triphenylene | 278 | C22H14 | 000215-58-7 | 96 |
| 2 | Picene | 278 | C22H14 | 000213-46-7 | 93 |
| 3 | Benzo[b]chrysene | 278 | C22H14 | 000214-17-5 | 86 |
| 4 | Dibenz[a,h]anthracene | 278 | C22H14 | 000053-70-3 | 86 |
| 5 | Vanadium, (.eta.5-2,4-cyclopenta... | 278 | C17H23V | 126948-97-8 | 64 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

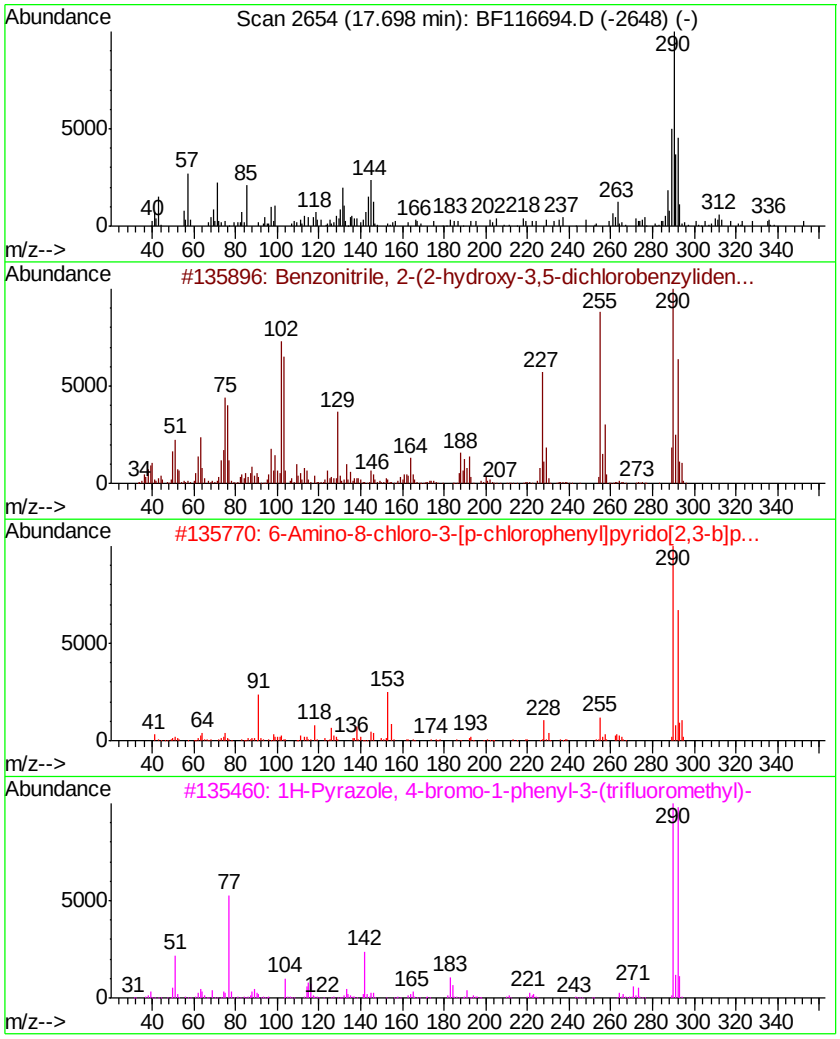
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

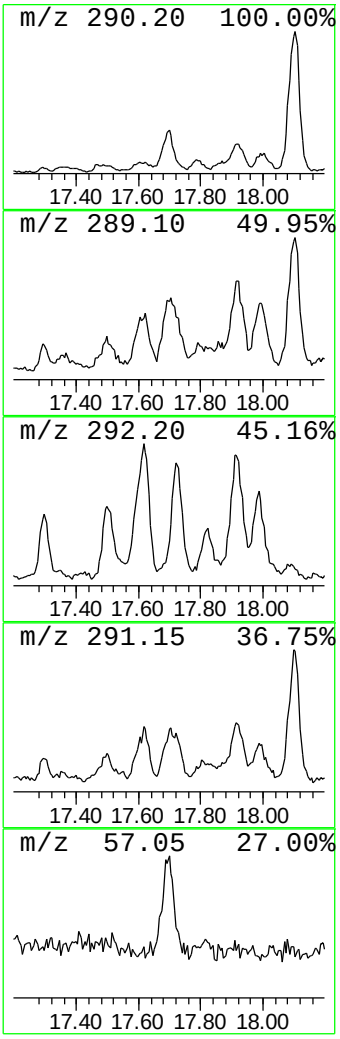
**********

Peak Number 18 Benzonitrile, 2-(2-hydroxy-... Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.70 | 5.69 ng | 220272 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzonitrile, 2-(2-hydroxy-3,5-d... | 290 | C14H8Cl2N2O | 316133-55-8 | 86 |
| 2 | 6-Amino-8-chloro-3-[p-chlorophen... | 290 | C13H8Cl2N4 | 028649-00-5 | 43 |
| 3 | 1H-Pyrazole, 4-bromo-1-phenyl-3-... | 290 | C10H6BrF3N2 | 1000327-39-4 | 43 |
| 4 | Benzo[ghi]perylene, 4-methyl- | 290 | C23H14 | 019224-38-5 | 38 |
| 5 | 3-Bromo-2,5,6-trimethoxybenzoic ... | 290 | C10H11BrO5 | 101460-22-4 | 38 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

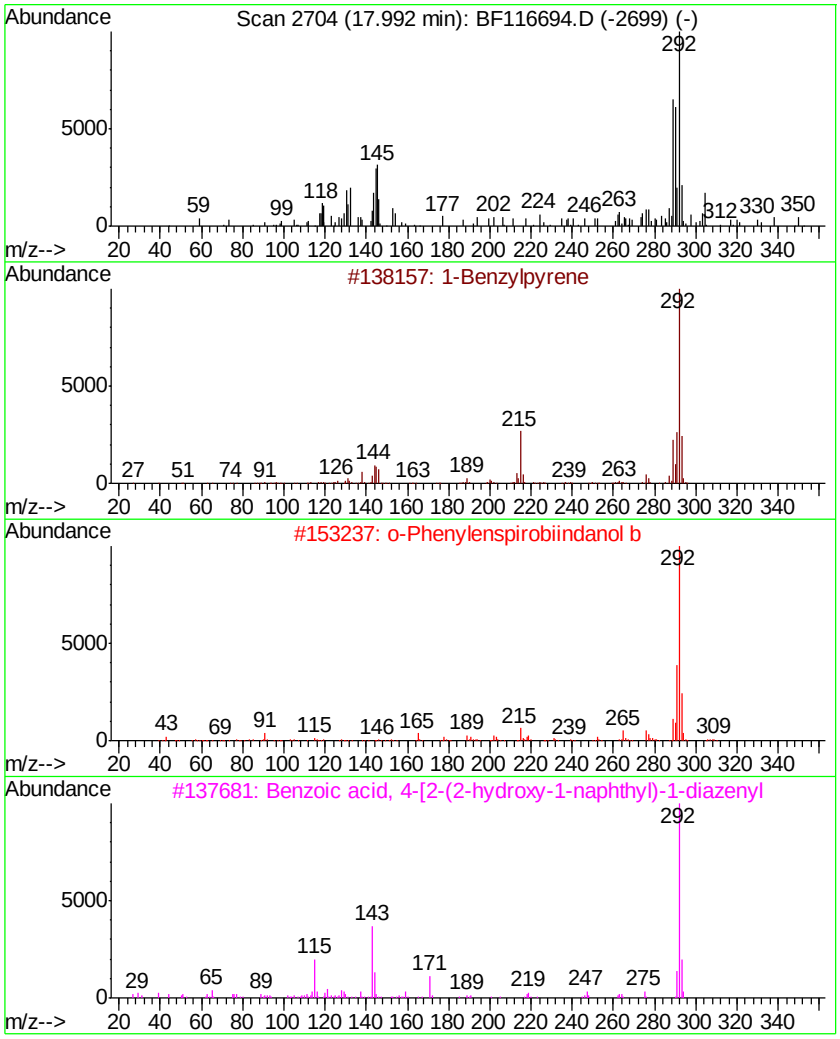
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

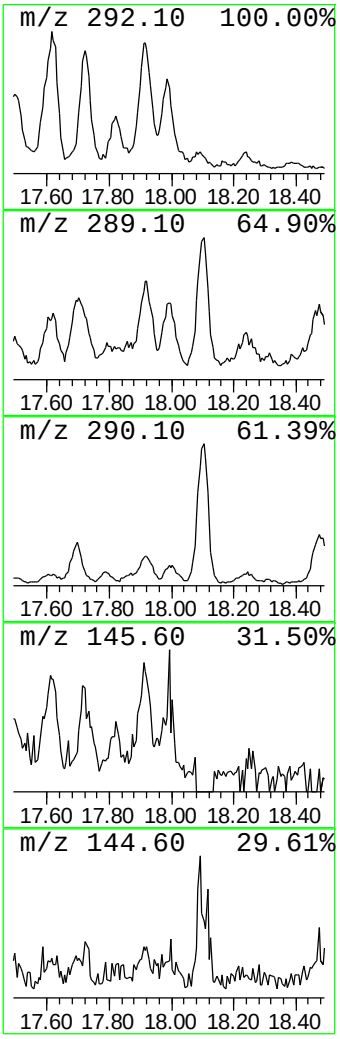
**********

Peak Number 19 unknown17.99 Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.99 | 4.24 ng | 164280 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Benzylpyrene | 292 | C23H16 | 042211-34-7 | 37 |
| 2 | o-Phenylenspirobiindanol b | 310 | C23H18O | 155919-26-9 | 35 |
| 3 | Benzoic acid, 4-[2-(2-hydroxy-1-... | 292 | C17H12N2O3 | 1000197-03-1 | 32 |
| 4 | Androstane-3,17-diol, (3.beta.,5... | 292 | C19H32O2 | 000571-20-0 | 30 |
| 5 | Aniline, p,p'-(m-phenylenedioxy)di- | 292 | C18H16N2O2 | 002479-46-1 | 27 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

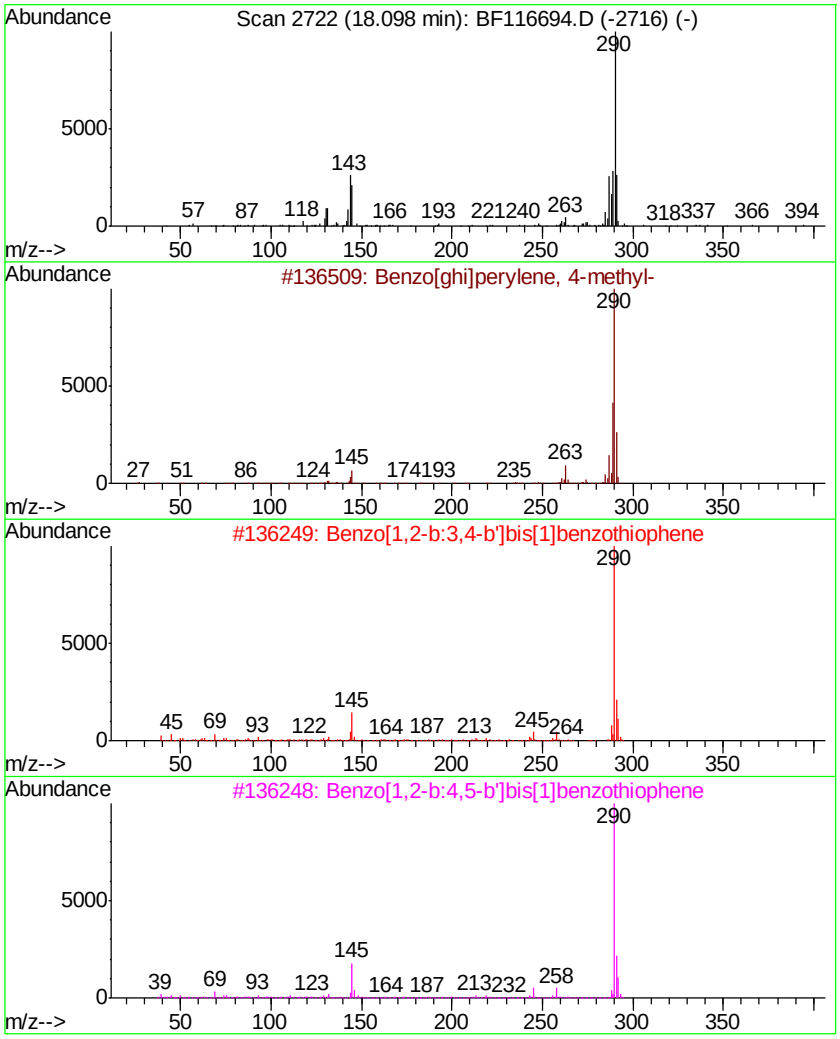
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

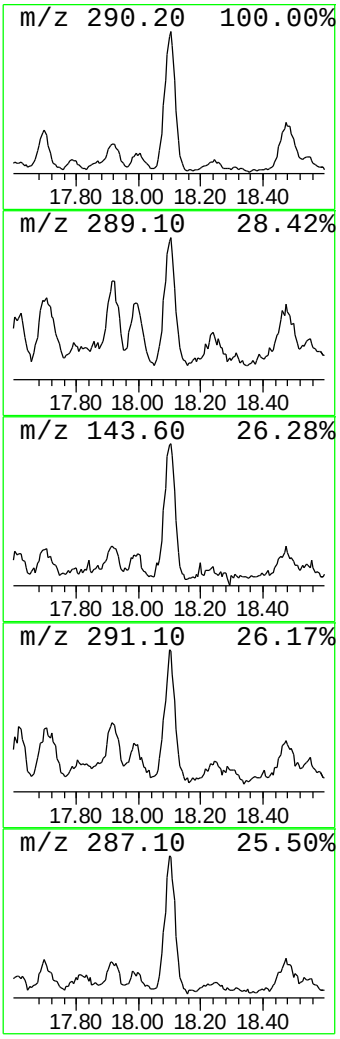
**********

Peak Number 20 Benzo[ghi]perylene, 4-methyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.10 | 11.48 ng | 444681 | Perylene-d12 | 15.45 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[ghi]perylene, 4-methyl- | 290 | C23H14 | 019224-38-5 | 74 |
| 2 | Benzo[1,2-b:3,4-b']bis[1]benzoth... | 290 | C18H10S2 | 000198-89-0 | 50 |
| 3 | Benzo[1,2-b:4,5-b']bis[1]benzoth... | 290 | C18H10S2 | 000241-34-9 | 47 |
| 4 | (4-Nitrophenyl)diphenylamine | 290 | C18H14N2O2 | 004316-57-8 | 38 |
| 5 | 10-Hexyl-9-anthracenecarbaldehyde | 290 | C21H22O | 165663-08-1 | 38 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF083119\
Data File : BF116694.D
Acq On : 31 Aug 2019 12:56
Operator : HP/JU
Sample : K4618-07 20X
Misc :
ALS Vial : 8 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-4-20-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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 |
| Chrysene, 1-methyl- | 14.39 | 14.9 | ng | 1701640 | 5 | 14.00 | 2287140 | 20.0 |
| Triphenylene, 2-m... | 14.42 | 7.8 | ng | 897677 | 5 | 14.00 | 2287140 | 20.0 |
| unknown14.47 | 14.47 | 5.7 | ng | 650761 | 5 | 14.00 | 2287140 | 20.0 |
| Benz[a]anthracene... | 14.72 | 5.9 | ng | 679067 | 5 | 14.00 | 2287140 | 20.0 |
| Benzo[c]phenanthr... | 14.77 | 13.3 | ng | 516624 | 6 | 15.45 | 774665 | 20.0 |
| Benz(a)anthracene... | 14.80 | 11.2 | ng | 433312 | 6 | 15.45 | 774665 | 20.0 |
| Benzo[c]phenanthr... | 14.87 | 4.6 | ng | 176459 | 6 | 15.45 | 774665 | 20.0 |
| 4H-Dibenz[a,kl]an... | 14.93 | 13.0 | ng | 504600 | 6 | 15.45 | 774665 | 20.0 |
| Tetradecane | 15.11 | 10.1 | ng | 389996 | 6 | 15.45 | 774665 | 20.0 |
| Benzo[e]pyrene | 15.33 | 25.9 | ng | 1002750 | 6 | 15.45 | 774665 | 20.0 |
| 11H-Indeno[2,1-a]... | 15.62 | 6.1 | ng | 236220 | 6 | 15.45 | 774665 | 20.0 |
| Perylene, 3-methyl- | 15.77 | 15.7 | ng | 607437 | 6 | 15.45 | 774665 | 20.0 |
| 10-Methylbenzo(a)... | 15.81 | 16.1 | ng | 625583 | 6 | 15.45 | 774665 | 20.0 |
| Benz[j]aceanthryl... | 15.89 | 8.1 | ng | 312821 | 6 | 15.45 | 774665 | 20.0 |
| 8H-Indeno[2,1-b]p... | 15.92 | 5.3 | ng | 203695 | 6 | 15.45 | 774665 | 20.0 |
| 7,10-Dimethylbenz... | 16.26 | 6.2 | ng | 240956 | 6 | 15.45 | 774665 | 20.0 |
| Benzo[b]triphenylene | 17.13 | 5.0 | ng | 194484 | 6 | 15.45 | 774665 | 20.0 |
| Benzonitrile, 2-(... | 17.70 | 5.7 | ng | 220272 | 6 | 15.45 | 774665 | 20.0 |
| unknown17.99 | 17.99 | 4.2 | ng | 164280 | 6 | 15.45 | 774665 | 20.0 |
| Benzo[ghi]perylen... | 18.10 | 11.5 | ng | 444681 | 6 | 15.45 | 774665 | 20.0 |