LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

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-BF022020.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 | 4.916 | 477 | 481 | 495 | rBV | 198090 | 295262 | 10.68% | 0.941% |
| 2 | 5.345 | 551 | 554 | 586 | rBV | 1339882 | 1915519 | 69.30% | 6.102% |
| 3 | 6.375 | 726 | 729 | 756 | rBV | 1626488 | 1932240 | 69.90% | 6.155% |
| 4 | 6.716 | 784 | 787 | 799 | rBV | 625231 | 576561 | 20.86% | 1.837% |
| 5 | 6.875 | 810 | 814 | 827 | rBV | 1735213 | 1735797 | 62.79% | 5.529% |
| 6 | 7.286 | 880 | 884 | 912 | rBV | 1207057 | 1325189 | 47.94% | 4.221% |
| 7 | 7.998 | 1002 | 1005 | 1023 | rVB | 773602 | 764552 | 27.66% | 2.435% |
| 8 | 9.075 | 1184 | 1188 | 1199 | rBV | 2823950 | 2525219 | 91.35% | 8.044% |
| 9 | 9.475 | 1253 | 1256 | 1262 | rVB | 175278 | 174538 | 6.31% | 0.556% |
| 10 | 9.751 | 1299 | 1303 | 1306 | rBV | 1074289 | 948344 | 34.31% | 3.021% |
| 11 | 9.786 | 1306 | 1309 | 1319 | rVB | 90368 | 96052 | 3.47% | 0.306% |
| 12 | 9.963 | 1334 | 1339 | 1343 | rBV2 | 34470 | 37319 | 1.35% | 0.119% |
| 13 | 10.304 | 1391 | 1397 | 1403 | rBV2 | 74972 | 83997 | 3.04% | 0.268% |
| 14 | 10.457 | 1421 | 1423 | 1426 | rVB | 29983 | 27701 | 1.00% | 0.088% |
| 15 | 10.545 | 1434 | 1438 | 1452 | rBV | 1987546 | 1871766 | 67.71% | 5.962% |
| 16 | 10.851 | 1483 | 1490 | 1493 | rBV3 | 37300 | 51891 | 1.88% | 0.165% |
| 17 | 10.933 | 1497 | 1504 | 1506 | rVB5 | 18011 | 28409 | 1.03% | 0.090% |
| 18 | 10.974 | 1506 | 1511 | 1513 | rBV2 | 36111 | 43222 | 1.56% | 0.138% |
| 19 | 11.092 | 1528 | 1531 | 1534 | rBV3 | 41672 | 47959 | 1.73% | 0.153% |
| 20 | 11.139 | 1536 | 1539 | 1544 | rVB | 40627 | 41929 | 1.52% | 0.134% |
| 21 | 11.239 | 1552 | 1556 | 1558 | rBV | 1125950 | 1069389 | 38.69% | 3.406% |
| 22 | 11.263 | 1558 | 1560 | 1564 | rVV | 879772 | 745942 | 26.99% | 2.376% |
| 23 | 11.316 | 1564 | 1569 | 1573 | rVV | 193422 | 187578 | 6.79% | 0.597% |
| 24 | 11.480 | 1592 | 1597 | 1603 | rBV3 | 74613 | 121527 | 4.40% | 0.387% |
| 25 | 11.527 | 1603 | 1605 | 1607 | rVB2 | 38195 | 28960 | 1.05% | 0.092% |
| 26 | 11.739 | 1638 | 1641 | 1643 | rBV | 146859 | 120974 | 4.38% | 0.385% |
| 27 | 11.769 | 1643 | 1646 | 1652 | rVV | 341685 | 359146 | 12.99% | 1.144% |
| 28 | 11.816 | 1652 | 1654 | 1657 | rVV | 84001 | 87284 | 3.16% | 0.278% |
| 29 | 11.851 | 1657 | 1660 | 1663 | rVV | 235002 | 264061 | 9.55% | 0.841% |
| 30 | 11.874 | 1663 | 1664 | 1670 | rVB | 106284 | 77696 | 2.81% | 0.247% |
| 31 | 12.033 | 1688 | 1691 | 1696 | rVB | 107787 | 101731 | 3.68% | 0.324% |
| 32 | 12.080 | 1696 | 1699 | 1702 | rBV | 69806 | 60692 | 2.20% | 0.193% |
| 33 | 12.180 | 1713 | 1716 | 1720 | rBV2 | 35147 | 36043 | 1.30% | 0.115% |
| 34 | 12.292 | 1731 | 1735 | 1737 | rBV | 92508 | 93602 | 3.39% | 0.298% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

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-BF022020.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 12.327 | 1737 | 1741 | 1744 | rVV3 | 53649 | 94172 | 3.41% | 0.300% |
| 36 | 12.392 | 1747 | 1752 | 1759 | rVB3 | 52684 | 89502 | 3.24% | 0.285% |
| 37 | 12.451 | 1759 | 1762 | 1767 | rBV | 1356953 | 1262716 | 45.68% | 4.022% |
| 38 | 12.521 | 1771 | 1774 | 1777 | rVV2 | 24903 | 34311 | 1.24% | 0.109% |
| 39 | 12.551 | 1777 | 1779 | 1785 | rVV3 | 46381 | 61688 | 2.23% | 0.196% |
| 40 | 12.604 | 1785 | 1788 | 1791 | rVV | 79164 | 77353 | 2.80% | 0.246% |
| 41 | 12.680 | 1795 | 1801 | 1807 | rVV | 1218318 | 1365256 | 49.39% | 4.349% |
| 42 | 12.733 | 1807 | 1810 | 1816 | rVB2 | 73541 | 87925 | 3.18% | 0.280% |
| 43 | 12.816 | 1820 | 1824 | 1828 | rVV | 2786773 | 2764259 | 100.00% | 8.805% |
| 44 | 12.857 | 1828 | 1831 | 1835 | rVB | 59149 | 67200 | 2.43% | 0.214% |
| 45 | 12.904 | 1835 | 1839 | 1844 | rBV2 | 81379 | 139837 | 5.06% | 0.445% |
| 46 | 12.992 | 1850 | 1854 | 1855 | rBV3 | 64356 | 80733 | 2.92% | 0.257% |
| 47 | 13.015 | 1855 | 1858 | 1861 | rVB | 166890 | 169414 | 6.13% | 0.540% |
| 48 | 13.080 | 1866 | 1869 | 1871 | rVV | 142693 | 132494 | 4.79% | 0.422% |
| 49 | 13.116 | 1871 | 1875 | 1883 | rVV3 | 137457 | 223055 | 8.07% | 0.711% |
| 50 | 13.174 | 1883 | 1885 | 1888 | rVB3 | 33580 | 31548 | 1.14% | 0.100% |
| 51 | 13.210 | 1888 | 1891 | 1894 | rBV | 71711 | 67717 | 2.45% | 0.216% |
| 52 | 13.245 | 1894 | 1897 | 1906 | rVB3 | 64135 | 96752 | 3.50% | 0.308% |
| 53 | 13.386 | 1918 | 1921 | 1923 | rBV2 | 41828 | 37918 | 1.37% | 0.121% |
| 54 | 13.533 | 1942 | 1946 | 1952 | rVB2 | 46502 | 75110 | 2.72% | 0.239% |
| 55 | 13.657 | 1964 | 1967 | 1970 | rVB2 | 92553 | 98443 | 3.56% | 0.314% |
| 56 | 13.686 | 1970 | 1972 | 1974 | rBV | 68569 | 64938 | 2.35% | 0.207% |
| 57 | 13.715 | 1974 | 1977 | 1981 | rBV3 | 160777 | 255893 | 9.26% | 0.815% |
| 58 | 13.804 | 1987 | 1992 | 1996 | rVB | 24646 | 39710 | 1.44% | 0.126% |
| 59 | 13.880 | 1998 | 2005 | 2008 | rBV2 | 1086877 | 1478093 | 53.47% | 4.708% |
| 60 | 13.910 | 2008 | 2010 | 2013 | rVB | 418602 | 373349 | 13.51% | 1.189% |
| 61 | 13.992 | 2021 | 2024 | 2028 | rVB | 85592 | 88898 | 3.22% | 0.283% |
| 62 | 14.027 | 2028 | 2030 | 2032 | rBV2 | 42890 | 45515 | 1.65% | 0.145% |
| 63 | 14.051 | 2032 | 2034 | 2038 | rVV3 | 43768 | 55700 | 2.02% | 0.177% |
| 64 | 14.092 | 2038 | 2041 | 2044 | rVV | 27548 | 47566 | 1.72% | 0.152% |
| 65 | 14.121 | 2044 | 2046 | 2049 | rVB | 40479 | 38714 | 1.40% | 0.123% |
| 66 | 14.274 | 2067 | 2072 | 2075 | rBV | 112166 | 150230 | 5.43% | 0.479% |
| 67 | 14.304 | 2075 | 2077 | 2080 | rVV | 43183 | 47495 | 1.72% | 0.151% |
| 68 | 14.333 | 2080 | 2082 | 2086 | rVV3 | 75704 | 107889 | 3.90% | 0.344% |
| 69 | 14.374 | 2086 | 2089 | 2091 | rVV2 | 53844 | 67503 | 2.44% | 0.215% |
| 70 | 14.398 | 2091 | 2093 | 2095 | rVV2 | 60045 | 68645 | 2.48% | 0.219% |
| 71 | 14.462 | 2099 | 2104 | 2107 | rVB2 | 28285 | 43265 | 1.57% | 0.138% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

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-BF022020.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 14.627 | 2130 | 2132 | 2135 | rVB | 64407 | 56133 | 2.03% | 0.179% |
| 73 | 14.692 | 2140 | 2143 | 2146 | rVB3 | 27907 | 35514 | 1.28% | 0.113% |
| 74 | 14.762 | 2151 | 2155 | 2160 | rBV3 | 56088 | 86907 | 3.14% | 0.277% |
| 75 | 14.862 | 2169 | 2172 | 2176 | rBV | 26576 | 34247 | 1.24% | 0.109% |
| 76 | 14.910 | 2176 | 2180 | 2183 | rVV | 385084 | 549318 | 19.87% | 1.750% |
| 77 | 14.939 | 2183 | 2185 | 2190 | rVB2 | 204679 | 236298 | 8.55% | 0.753% |
| 78 | 15.015 | 2195 | 2198 | 2203 | rVB | 87860 | 99531 | 3.60% | 0.317% |
| 79 | 15.198 | 2225 | 2229 | 2233 | rBV | 185547 | 225550 | 8.16% | 0.718% |
| 80 | 15.257 | 2236 | 2239 | 2243 | rBV | 326498 | 372325 | 13.47% | 1.186% |
| 81 | 15.309 | 2243 | 2248 | 2253 | rBV | 587400 | 730381 | 26.42% | 2.326% |
| 82 | 15.486 | 2274 | 2278 | 2282 | rBV3 | 34811 | 56582 | 2.05% | 0.180% |
| 83 | 15.698 | 2310 | 2314 | 2318 | rBV | 58062 | 77393 | 2.80% | 0.247% |
| 84 | 15.792 | 2325 | 2330 | 2338 | rBV5 | 32618 | 90888 | 3.29% | 0.290% |
| 85 | 15.862 | 2339 | 2342 | 2351 | rVB5 | 32446 | 53048 | 1.92% | 0.169% |
| 86 | 16.162 | 2389 | 2393 | 2397 | rBV3 | 31819 | 53407 | 1.93% | 0.170% |
| 87 | 16.533 | 2452 | 2456 | 2460 | rVB3 | 22010 | 32563 | 1.18% | 0.104% |
| 88 | 16.733 | 2482 | 2490 | 2498 | rVB | 138031 | 301301 | 10.90% | 0.960% |
| 89 | 16.892 | 2513 | 2517 | 2523 | rVV2 | 29850 | 53013 | 1.92% | 0.169% |
| 90 | 16.945 | 2523 | 2526 | 2534 | rVB2 | 29682 | 55049 | 1.99% | 0.175% |
| 91 | 17.156 | 2557 | 2562 | 2572 | rBV | 104551 | 225833 | 8.17% | 0.719% |
| 92 | 17.415 | 2602 | 2606 | 2613 | rVB | 31670 | 59806 | 2.16% | 0.191% |

Sum of corrected areas: 31393984

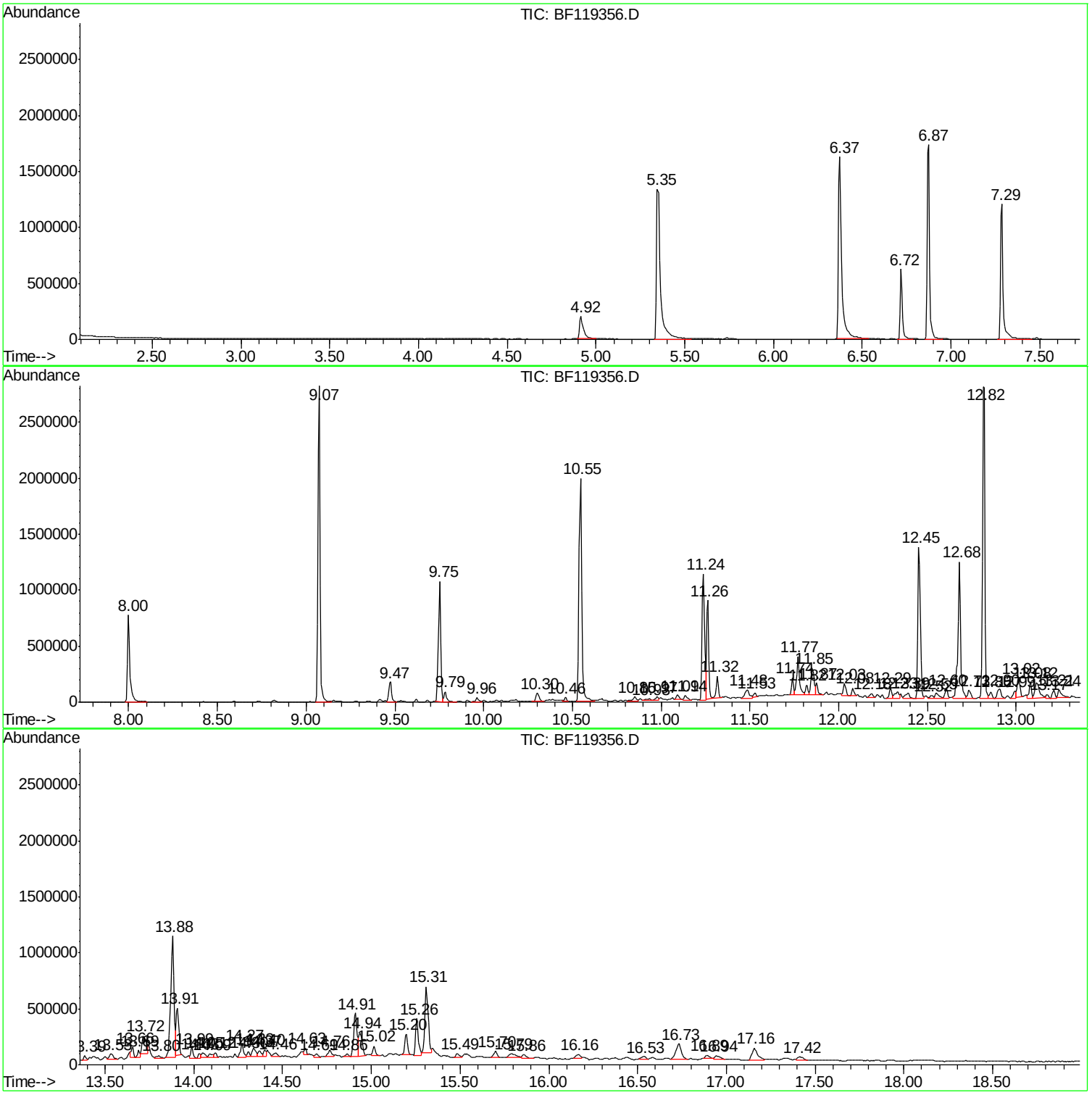
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.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\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

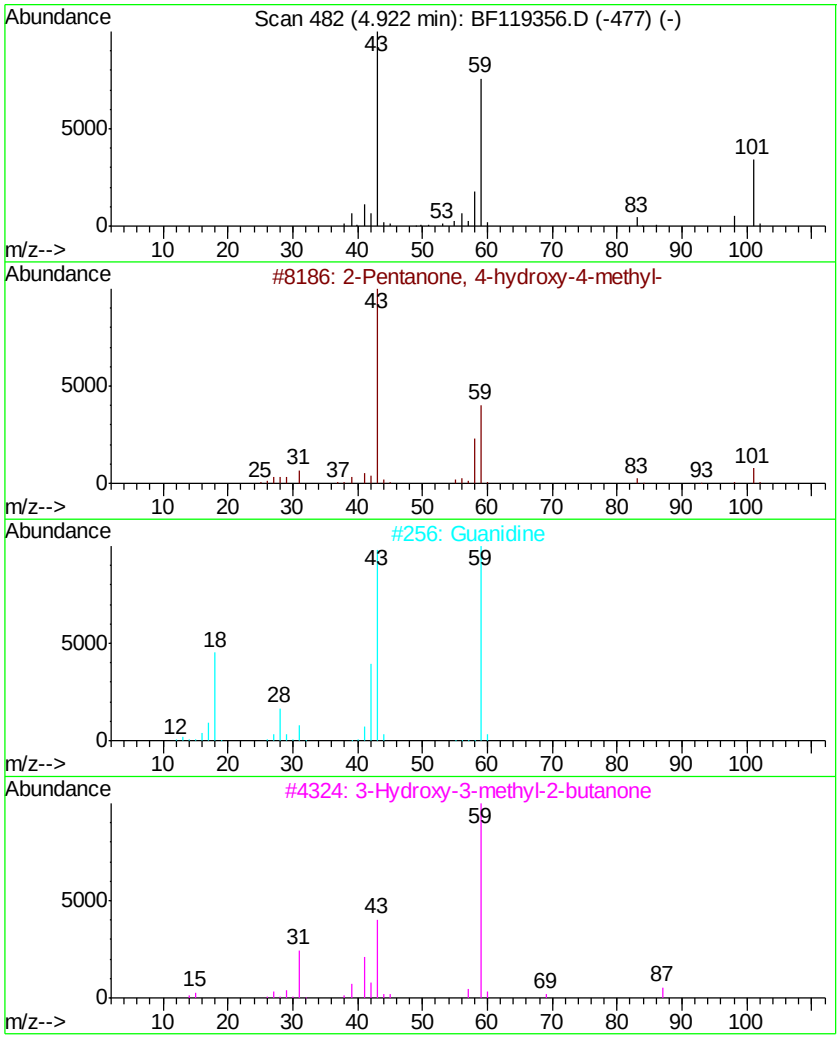
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

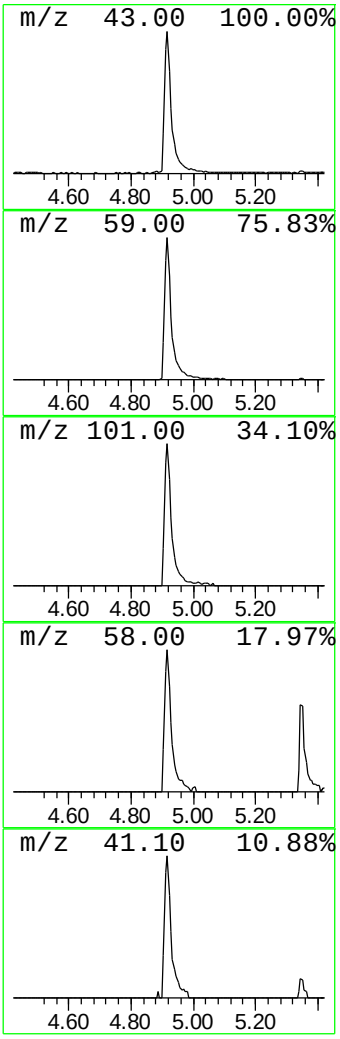
**********

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.92 | 10.24 ng | 295262 | 1,4-Dichlorobenzene-d4 | 6.72 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 56 |
| 2 | Guanidine | 59 | CH5N3 | 000113-00-8 | 50 |
| 3 | 3-Hydroxy-3-methyl-2-butanone | 102 | C5H10O2 | 000115-22-0 | 38 |
| 4 | (+-)-4-Amino-4,5-dihydro-2(3H)-f... | 101 | C4H7NO2 | 016504-58-8 | 35 |
| 5 | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 28 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

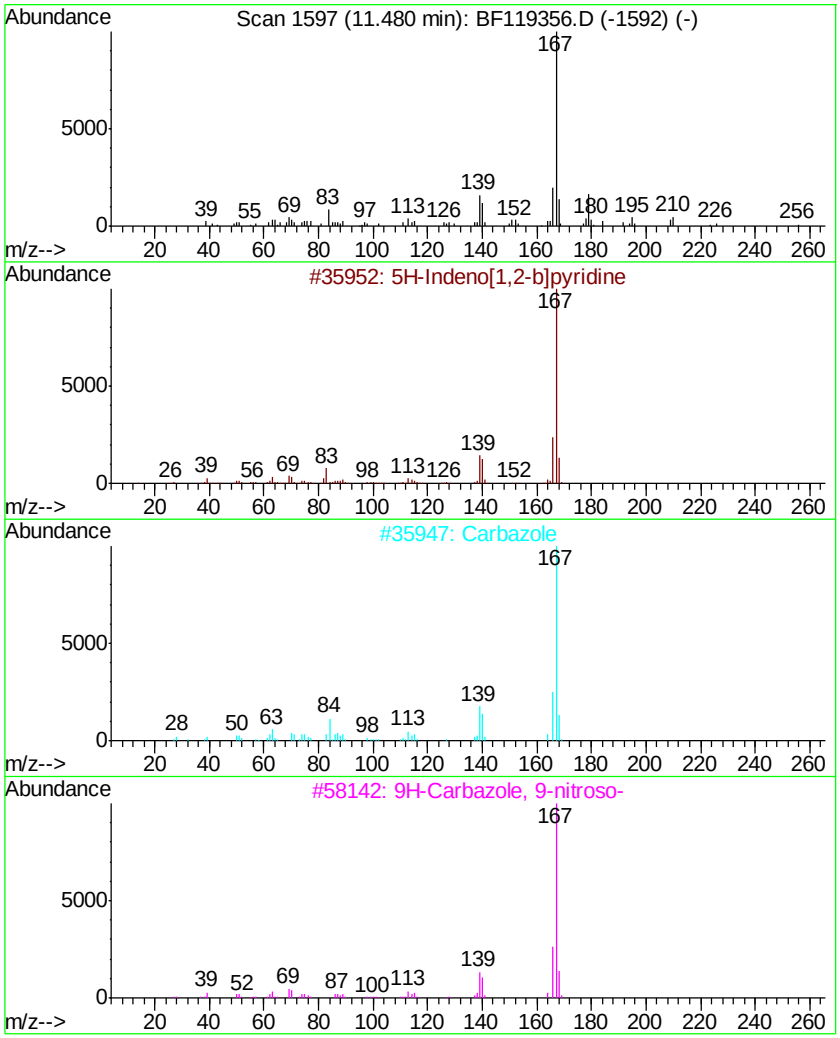
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

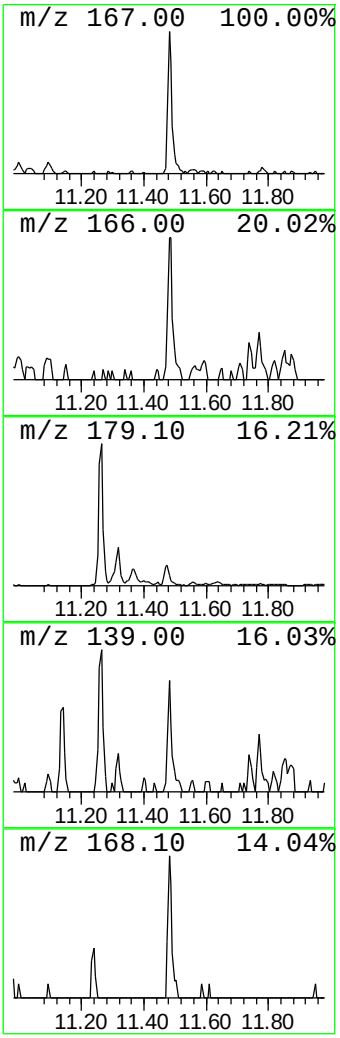
**********

Peak Number 3 5H-Indeno[1,2-b]pyridine Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.48 | 2.27 ng | 121527 | Phenanthrene-d10 | 11.24 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 5H-Indeno[1,2-b]pyridine | 167 | C12H9N | 000244-99-5 | 94 |
| 2 | Carbazole | 167 | C12H9N | 000086-74-8 | 90 |
| 3 | 9H-Carbazole, 9-nitroso- | 196 | C12H8N2O | 002788-23-0 | 74 |
| 4 | 9H-Carbazole-9-methanol | 197 | C13H11NO | 002409-36-1 | 74 |
| 5 | 2-Azafluorene | 167 | C12H9N | 000244-40-6 | 68 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

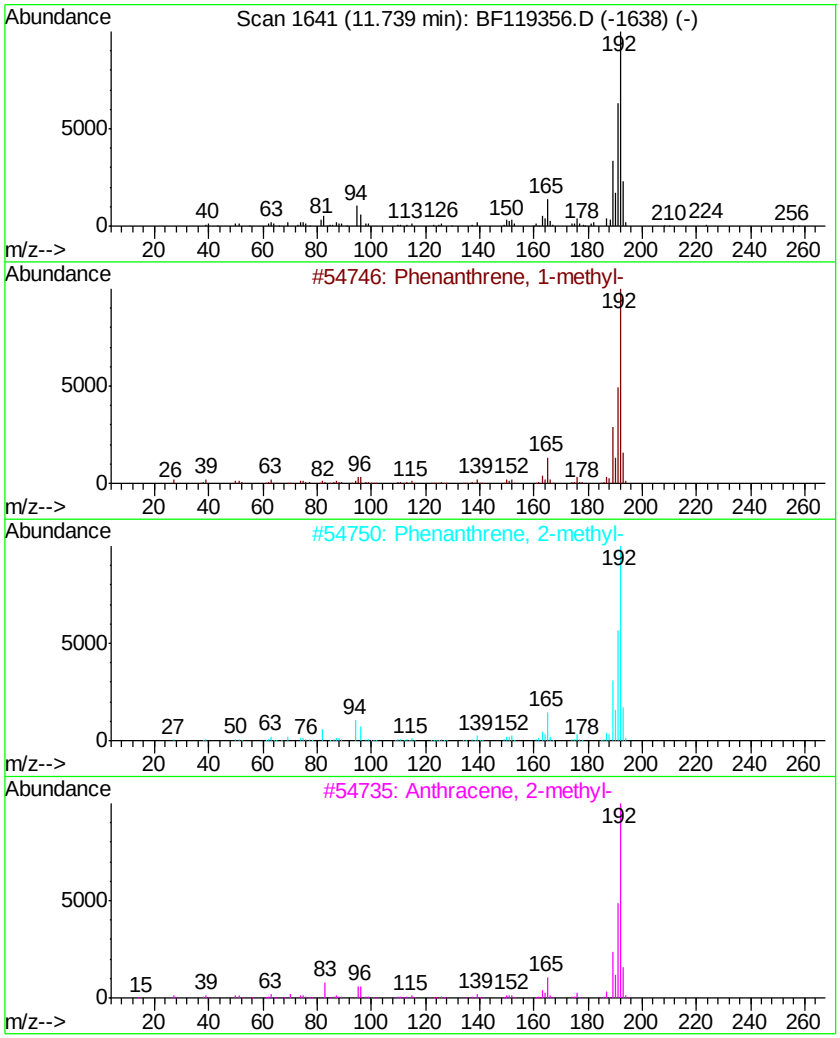
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

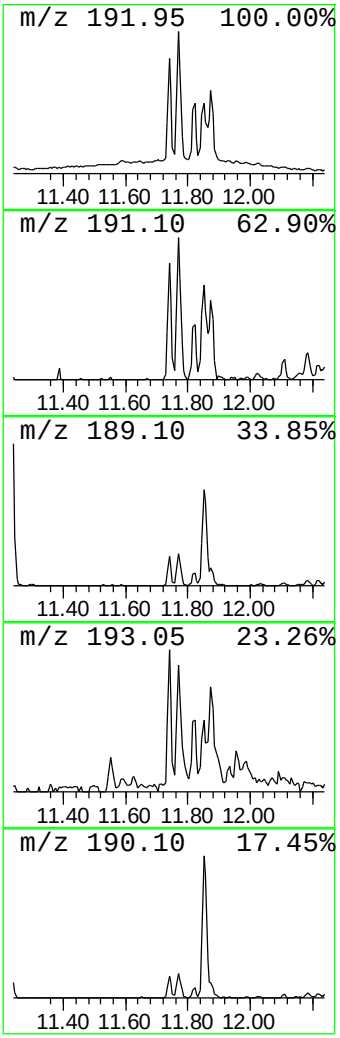
**********

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.74 | 2.26 ng | 120974 | Phenanthrene-d10 | 11.24 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 97 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 96 |
| 3 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 93 |
| 4 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 90 |
| 5 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 90 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

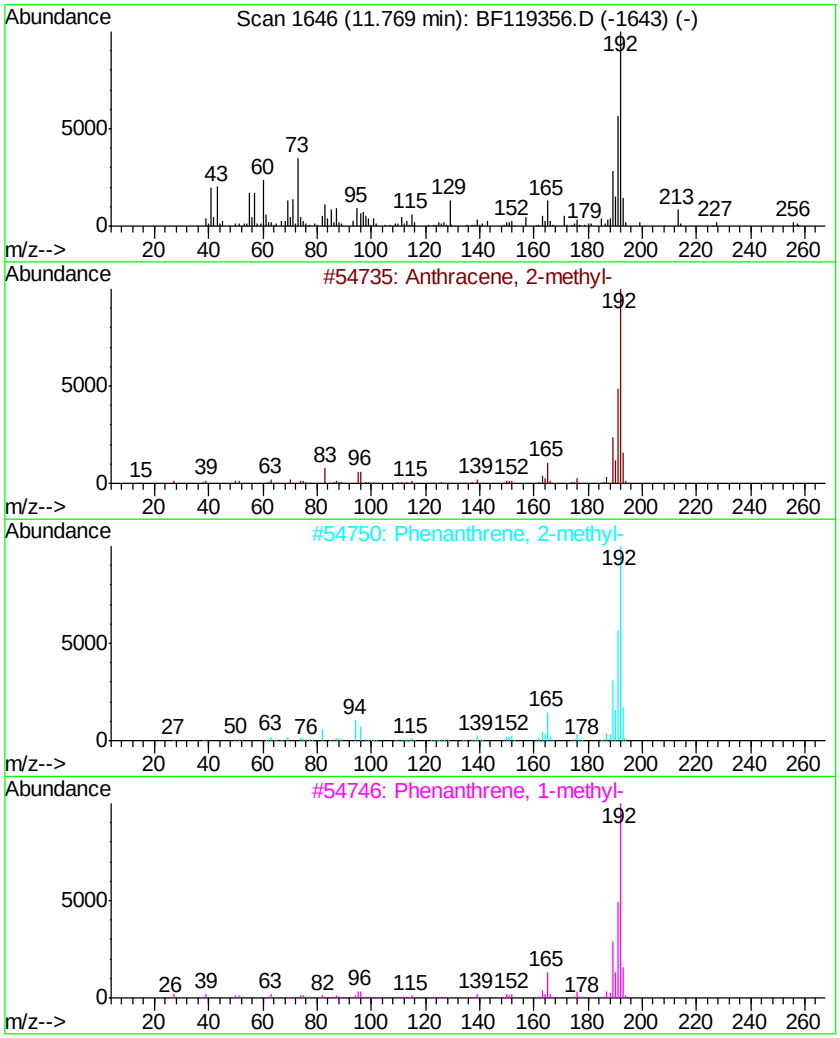
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

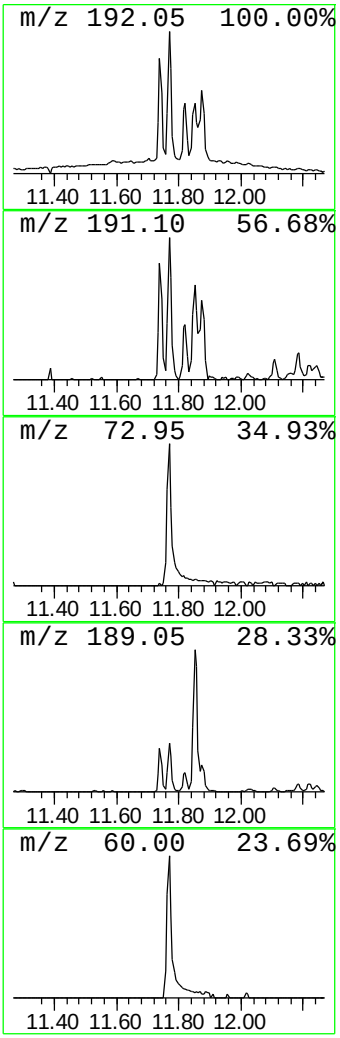
**********

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.77 | 6.72 ng | 359146 | Phenanthrene-d10 | 11.24 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 97 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 97 |
| 3 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 97 |
| 4 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 96 |
| 5 | Phenanthrene, 4-methyl- | 192 | C15H12 | 000832-64-4 | 91 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

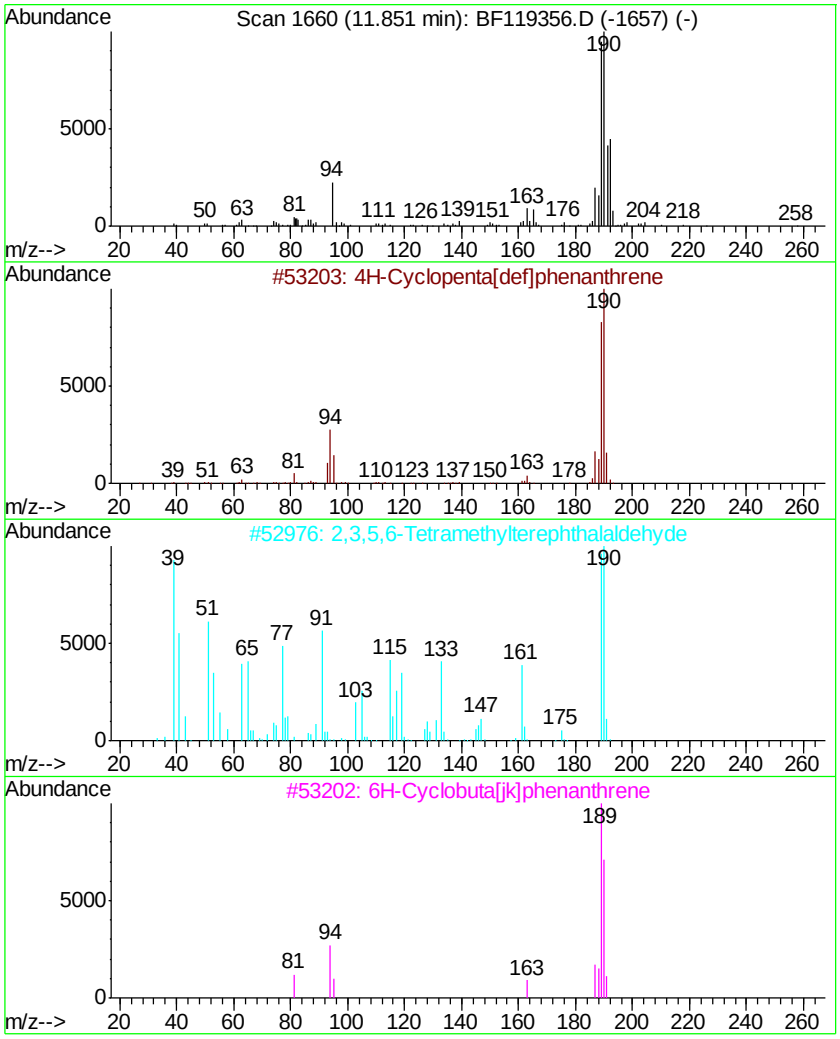
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

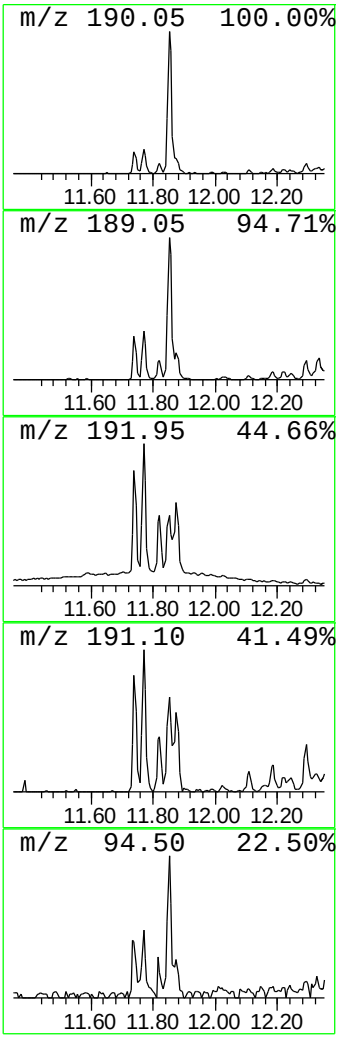
**********

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.85 | 4.94 ng | 264061 | Phenanthrene-d10 | 11.24 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 76 |
| 2 | 2,3,5,6-Tetramethylterephthalald... | 190 | C12H14O2 | 007072-01-7 | 43 |
| 3 | 6H-Cyclobuta[jk]phenanthrene | 190 | C15H10 | 083469-43-6 | 42 |
| 4 | 2,2'-Bis(4,5-dimethylimidazole) | 190 | C10H14N4 | 069286-06-2 | 40 |
| 5 | Carbonic acid, ethyl 2-formyl-4,... | 262 | C10H8Cl2O4 | 1000331-34-4 | 33 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

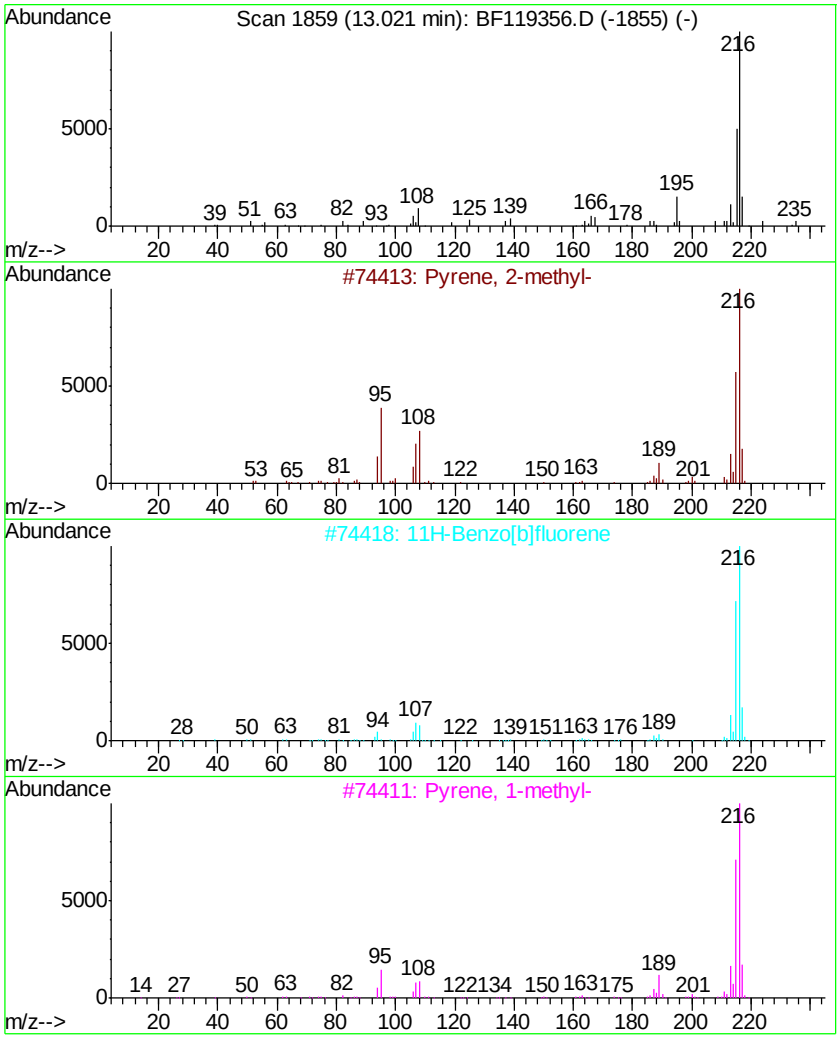
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

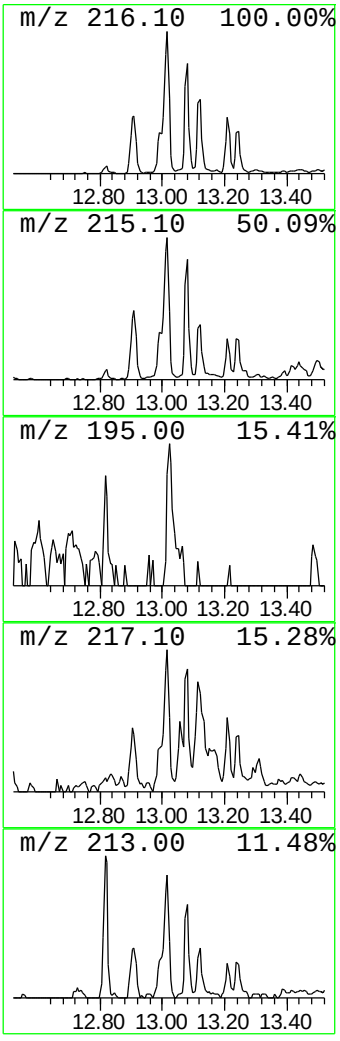
**********

Peak Number 7 Pyrene, 2-methyl- Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.02 | 2.29 ng | 169414 | Chrysene-d12 | 13.88 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 93 |
| 2 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 90 |
| 3 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 83 |
| 4 | 7H-Benzo[c]fluorene | 216 | C17H12 | 000205-12-9 | 74 |
| 5 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 72 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

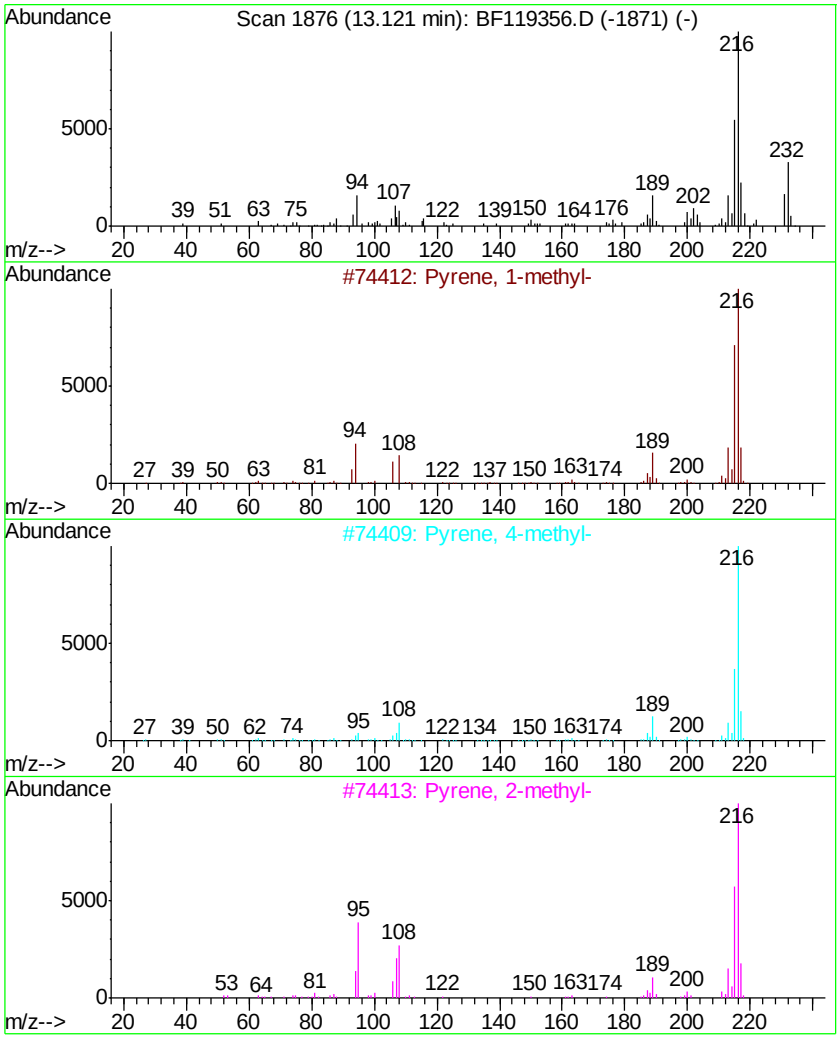
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

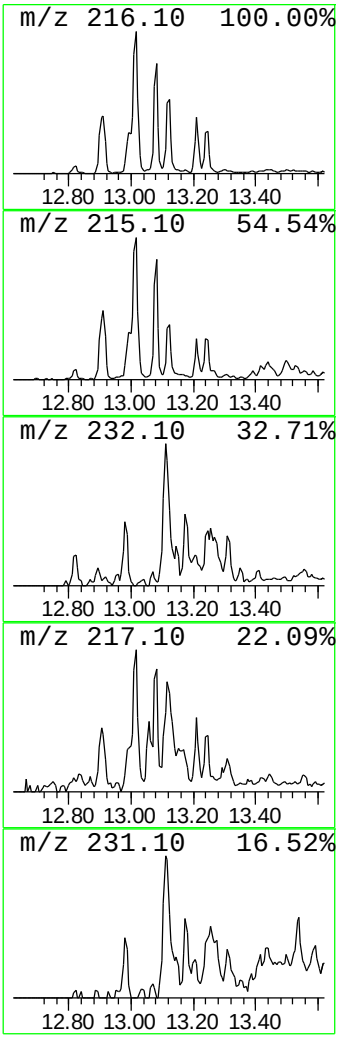
**********

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.12 | 3.02 ng | 223055 | Chrysene-d12 | 13.88 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 92 |
| 2 | Pyrene, 4-methyl- | 216 | C17H12 | 003353-12-6 | 55 |
| 3 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 49 |
| 4 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 46 |
| 5 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 43 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

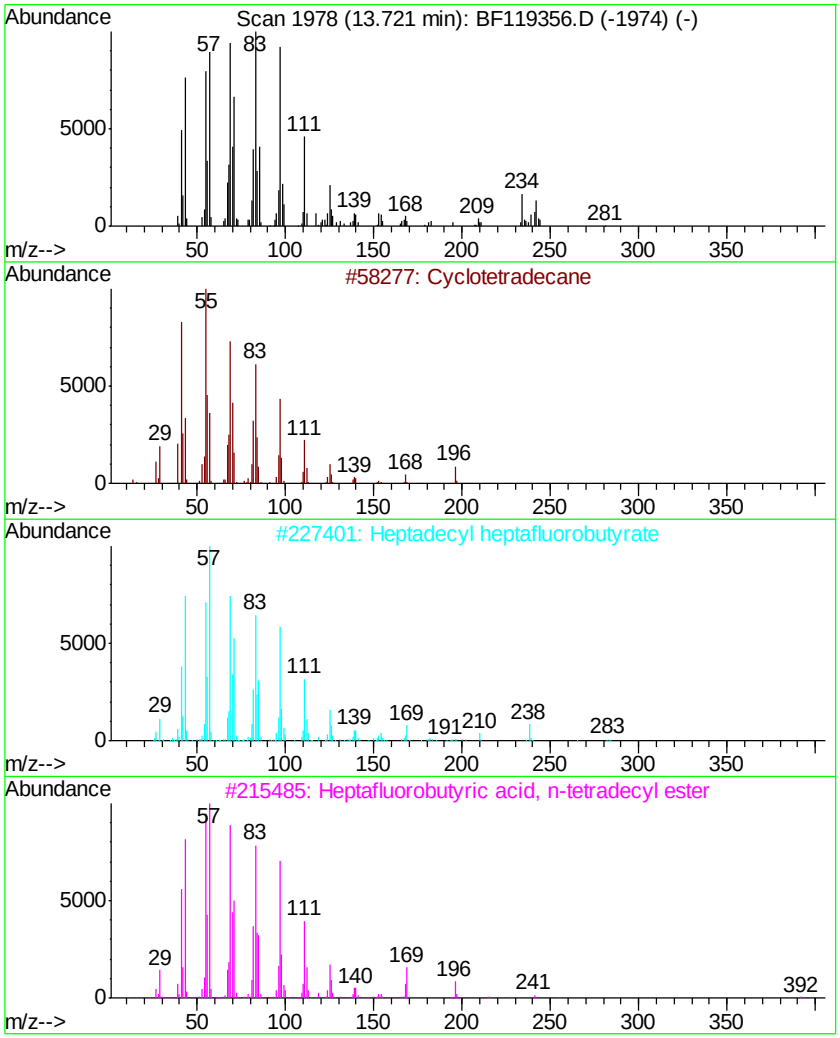
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

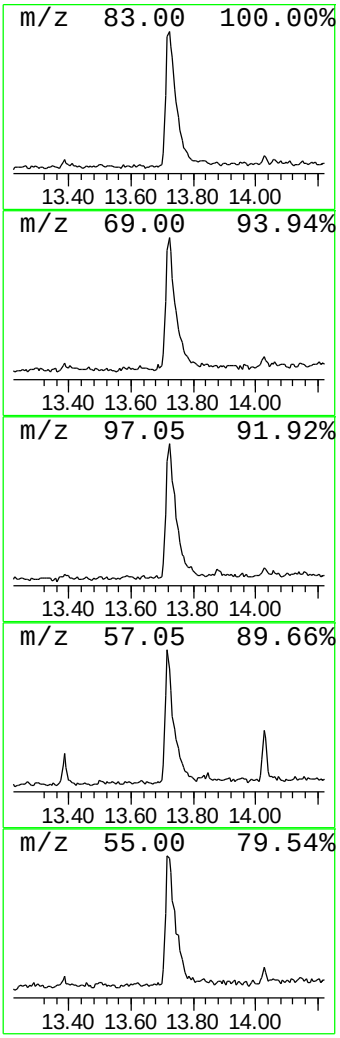
**********

Peak Number 9 Cyclotetradecane Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.72 | 3.46 ng | 255893 | Chrysene-d12 | 13.88 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclotetradecane | 196 | C14H28 | 000295-17-0 | 90 |
| 2 | Heptadecyl heptafluorobutyrate | 452 | C21H35F7O2 | 959085-66-6 | 90 |
| 3 | Heptafluorobutyric acid, n-tetra... | 410 | C18H29F7O2 | 007365-36-8 | 90 |
| 4 | Pentafluoropropionic acid, tetra... | 360 | C17H29F5O2 | 006222-06-6 | 90 |
| 5 | n-Nonadecanol-1 | 284 | C19H40O | 001454-84-8 | 87 |



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Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

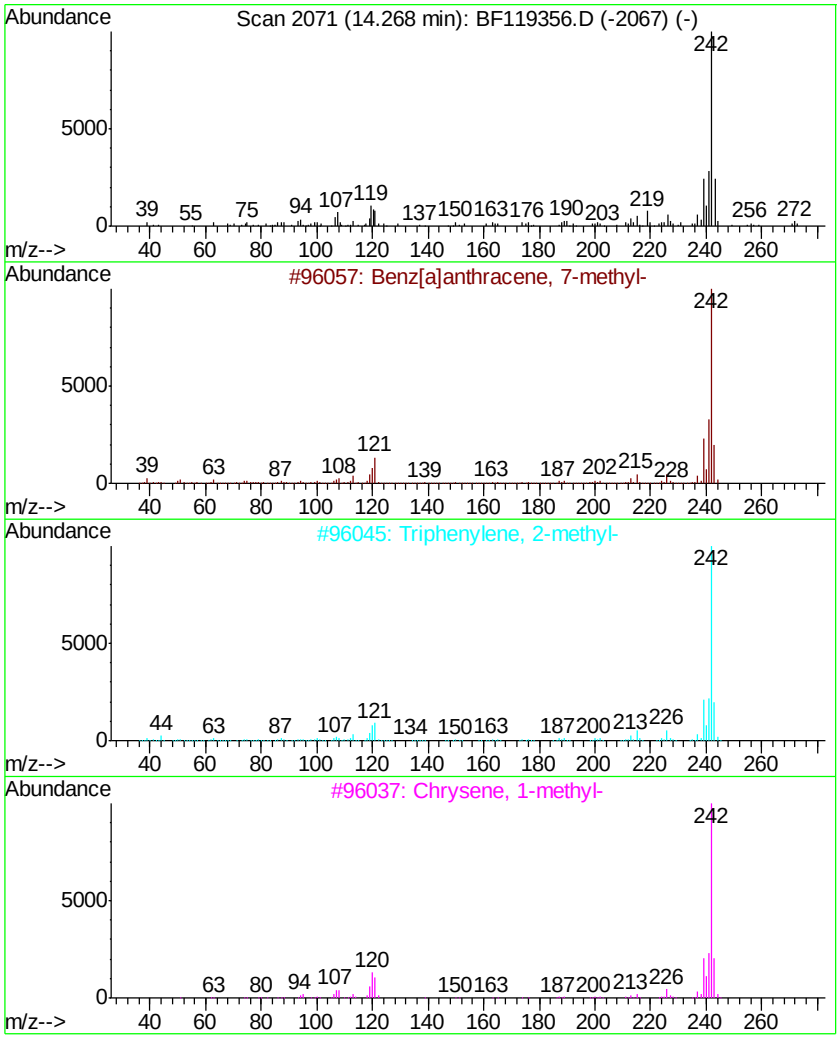
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

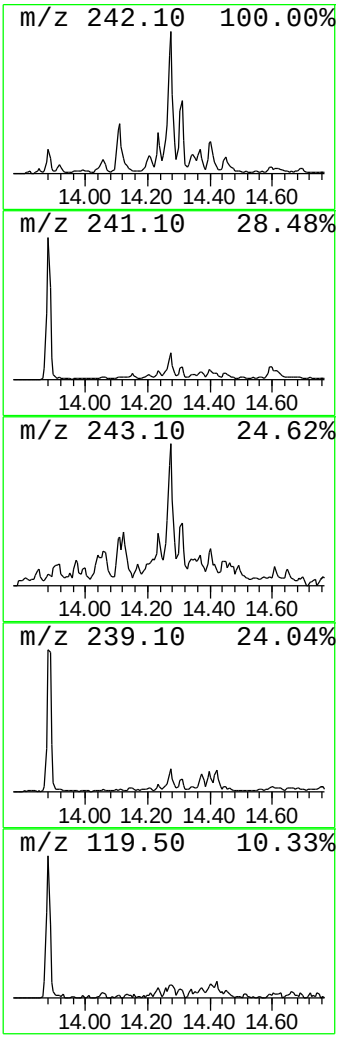
**********

Peak Number 10 Benz[a]anthracene, 7-methyl- Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.27 | 2.03 ng | 150230 | Chrysene-d12 | 13.88 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benz[a]anthracene, 7-methyl- | 242 | C19H14 | 002541-69-7 | 90 |
| 2 | Triphenylene, 2-methyl- | 242 | C19H14 | 001705-84-6 | 90 |
| 3 | Chrysene, 1-methyl- | 242 | C19H14 | 003351-28-8 | 90 |
| 4 | Benz[a]anthracene, 8-methyl- | 242 | C19H14 | 002381-31-9 | 90 |
| 5 | Chrysene, 3-methyl- | 242 | C19H14 | 003351-31-3 | 89 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

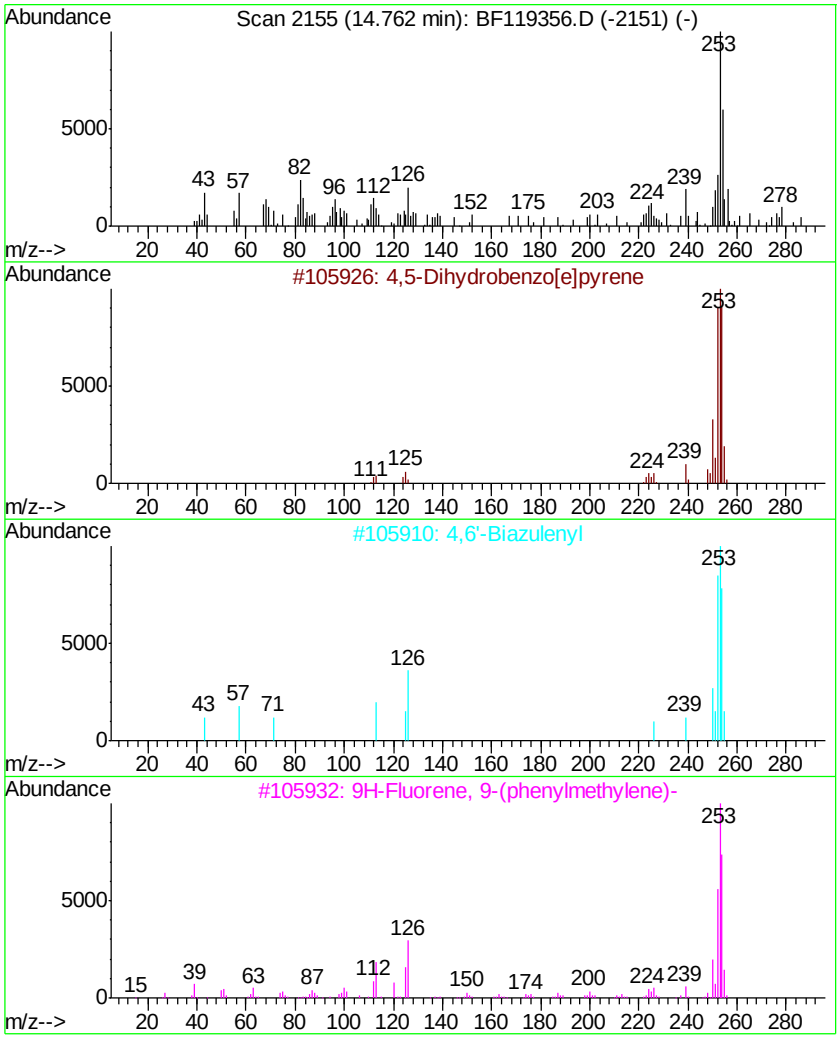
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

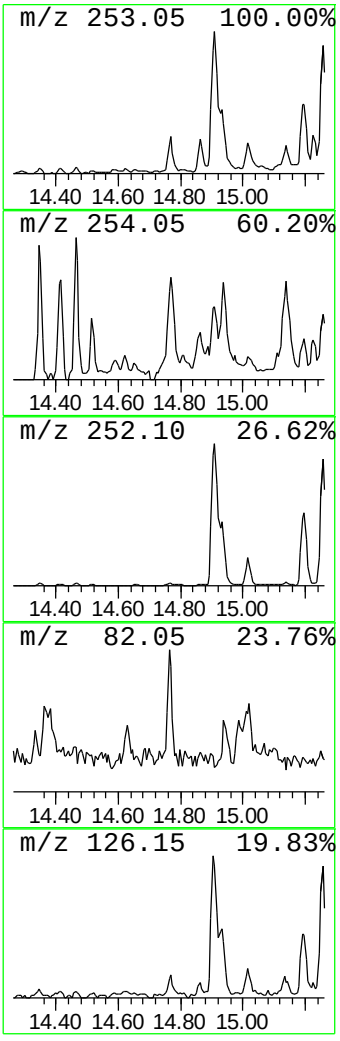
**********

Peak Number 11 4,5-Dihydrobenzo[e]pyrene Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.76 | 2.38 ng | 86907 | Perylene-d12 | 15.31 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4,5-Dihydrobenzo[e]pyrene | 254 | C20H14 | 095676-42-9 | 70 |
| 2 | 4,6'-Biazulenyl | 254 | C20H14 | 094154-49-1 | 62 |
| 3 | 9H-Fluorene, 9-(phenylmethylene)- | 254 | C20H14 | 001836-87-9 | 58 |
| 4 | 7,12-Dihydrobenzo[k]fluoranthene | 254 | C20H14 | 1000080-17-7 | 52 |
| 5 | Pyrido[2,3-g]indole, 5-methoxy-2... | 254 | C16H18N2O | 201991-45-9 | 52 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

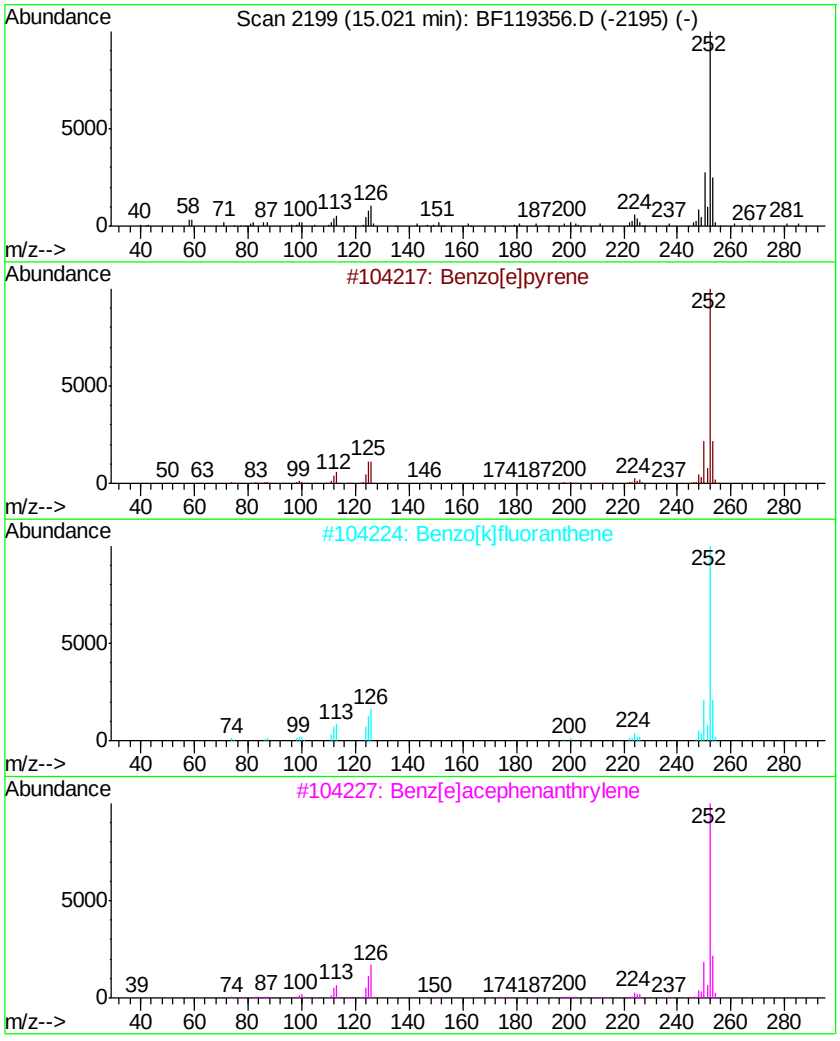
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

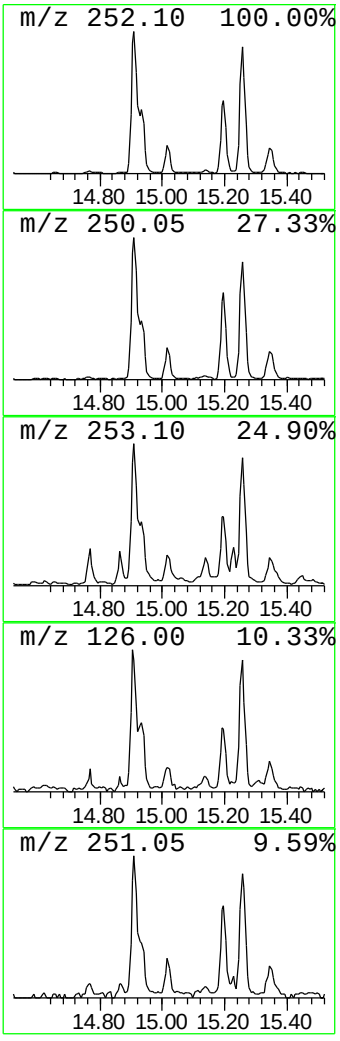
**********

Peak Number 12 Benzo[e]pyrene Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.02 | 2.73 ng | 99531 | Perylene-d12 | 15.31 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 94 |
| 2 | Benzo[k]fluoranthene | 252 | C20H12 | 000207-08-9 | 94 |
| 3 | Benz[e]acephenanthrylene | 252 | C20H12 | 000205-99-2 | 90 |
| 4 | Perylene | 252 | C20H12 | 000198-55-0 | 90 |
| 5 | Benzo[a]pyrene | 252 | C20H12 | 000050-32-8 | 81 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

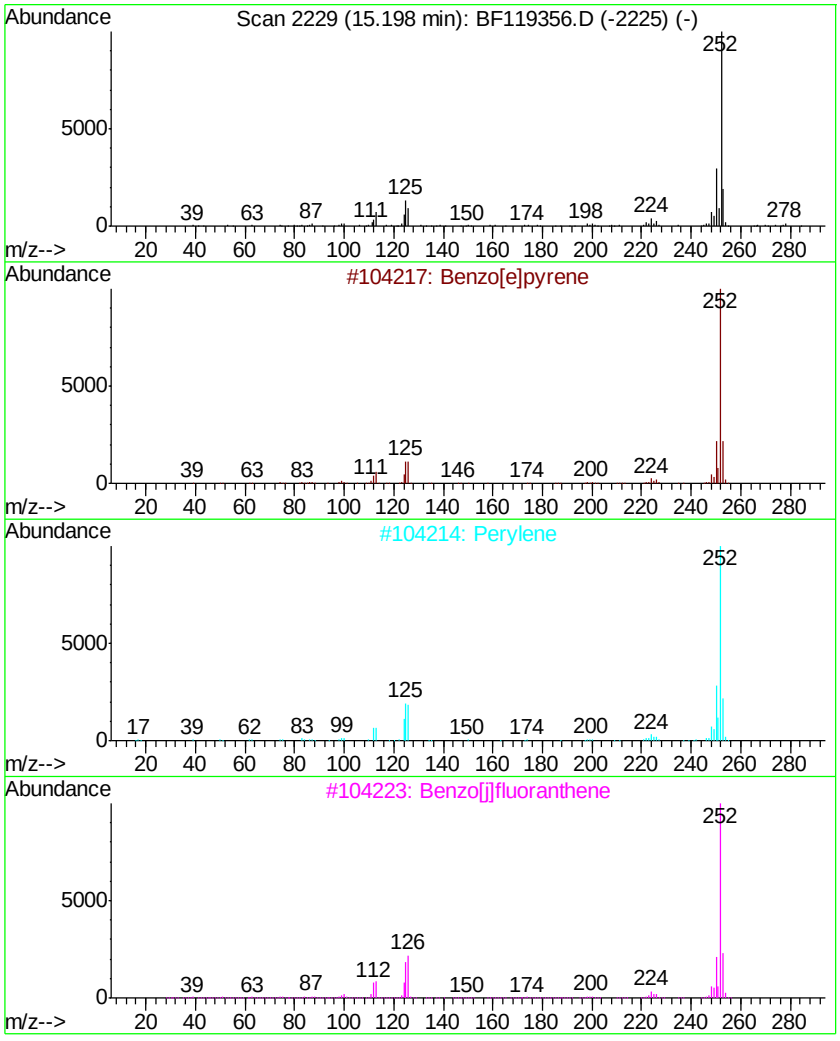
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

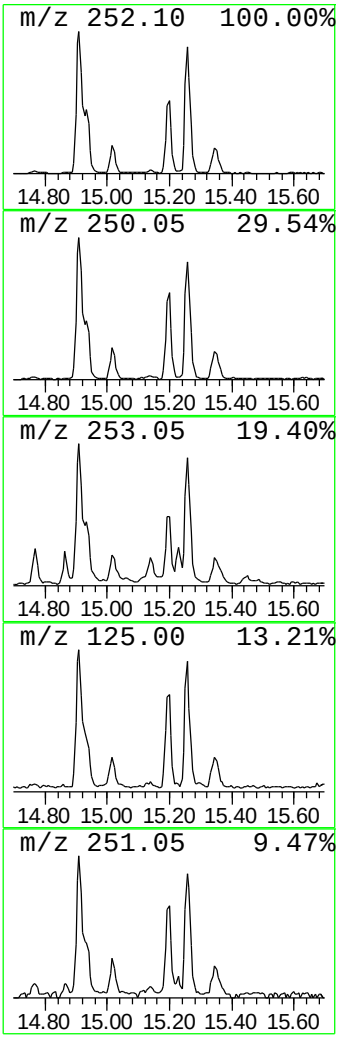
**********

Peak Number 13 Perylene Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.20 | 6.18 ng | 225550 | Perylene-d12 | 15.31 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 98 |
| 2 | Perylene | 252 | C20H12 | 000198-55-0 | 96 |
| 3 | Benzo[j]fluoranthene | 252 | C20H12 | 000205-82-3 | 94 |
| 4 | Benz[e]acephenanthrylene | 252 | C20H12 | 000205-99-2 | 90 |
| 5 | Benzo[k]fluoranthene | 252 | C20H12 | 000207-08-9 | 89 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

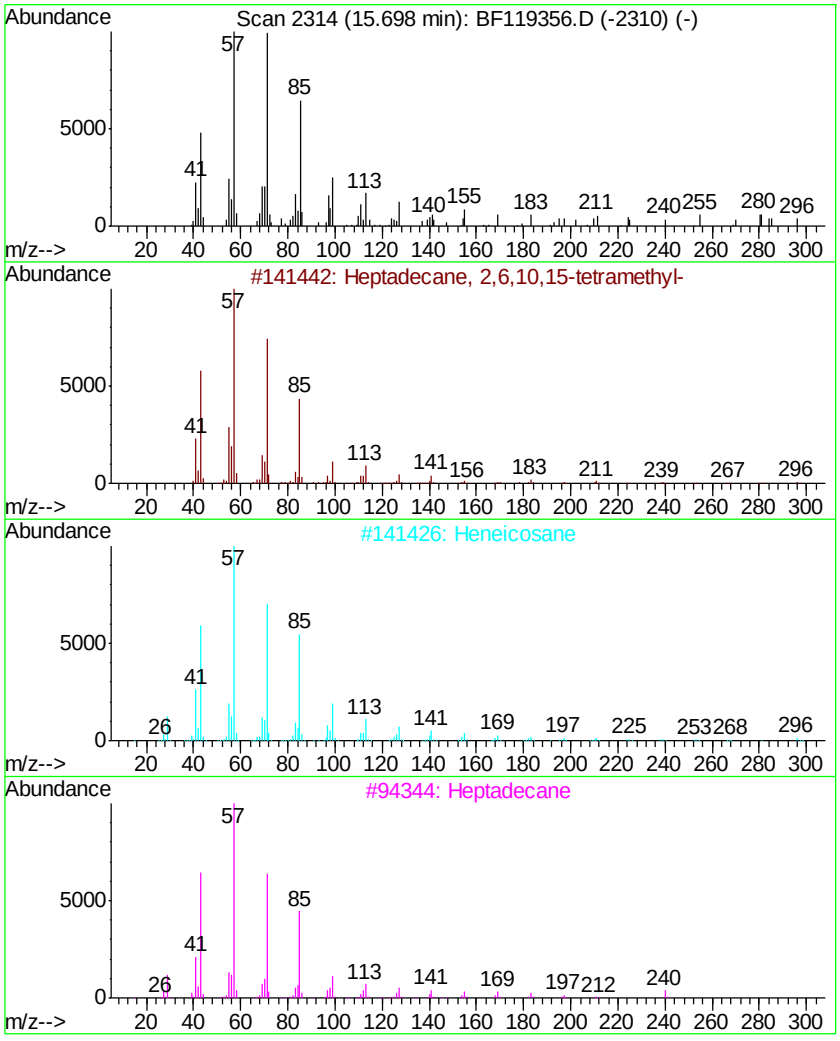
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

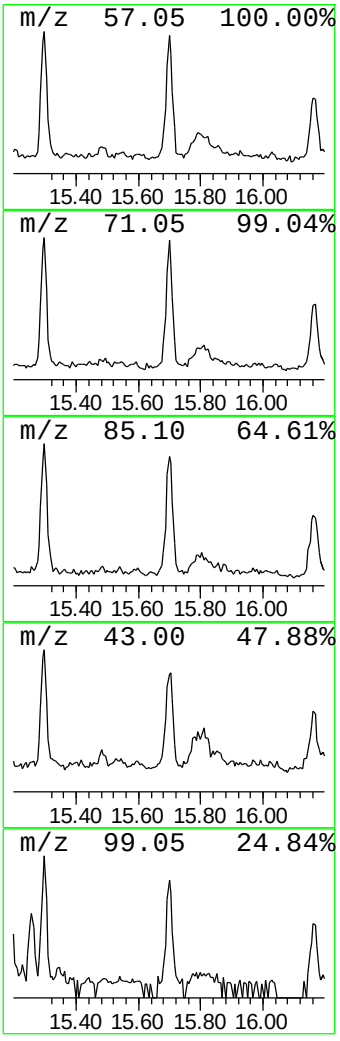
**********

Peak Number 14 Heptadecane, 2,6,10,15-tetr... Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.70 | 2.12 ng | 77393 | Perylene-d12 | 15.31 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44 | 054833-48-6 | 91 |
| 2 | Heneicosane | 296 | C21H44 | 000629-94-7 | 91 |
| 3 | Heptadecane | 240 | C17H36 | 000629-78-7 | 90 |
| 4 | Hexacosane | 366 | C26H54 | 000630-01-3 | 87 |
| 5 | Pentacosane | 352 | C25H52 | 000629-99-2 | 83 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.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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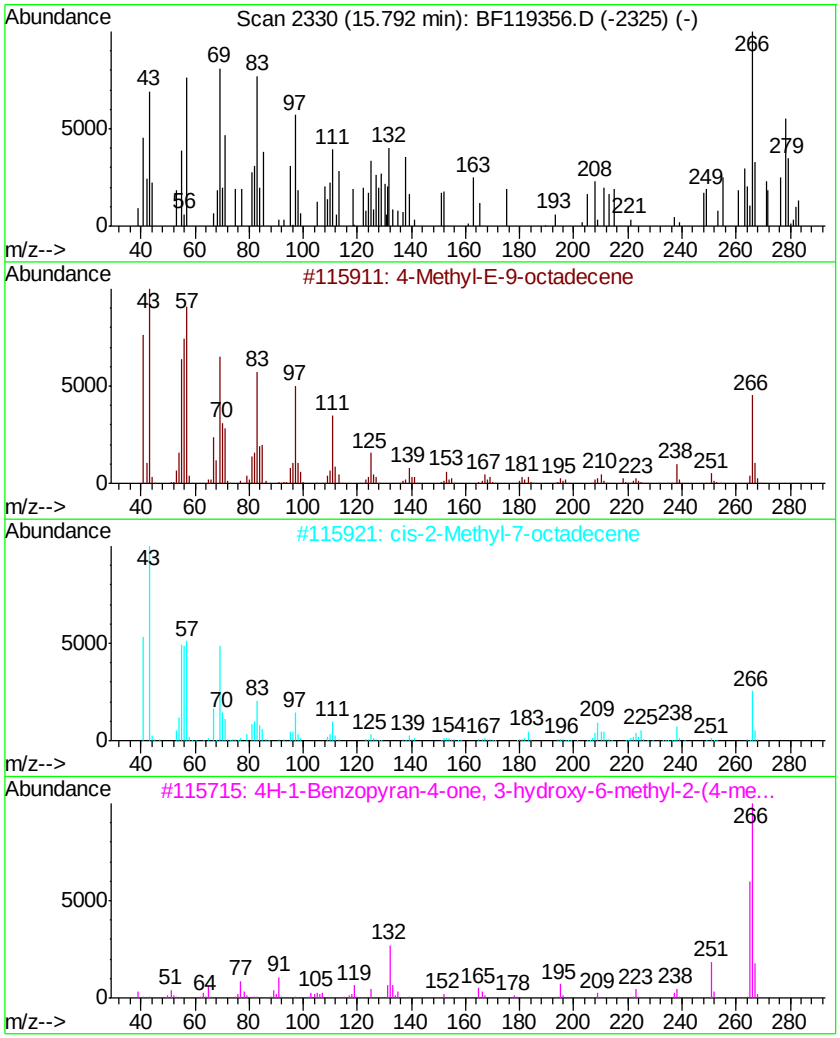
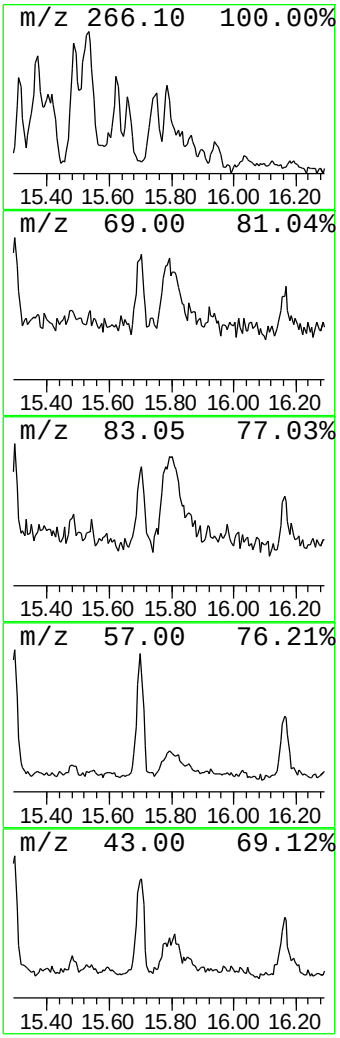
Peak Number 15 unknown15.79 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.79 | 2.49 ng | 90888 | Perylene-d12 | 15.31 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Methyl-E-9-octadecene | 266 | C19H38 | 1000130-87-1 | 35 |
| 2 | cis-2-Methyl-7-octadecene | 266 | C19H38 | 035354-39-3 | 32 |
| 3 | 4H-1-Benzopyran-4-one, 3-hydroxy... | 266 | C17H14O3 | 078396-37-9 | 22 |
| 4 | Imidazole-5-methanol, 2-ethyl-1-... | 266 | C17H18N2O | 309731-07-5 | 12 |
| 5 | Trifluoroacetoacetyl-2-naphthalene | 266 | C14H9F3O2 | 000893-33-4 | 11 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031120\
Data File : BF119356.D
Acq On : 11 Mar 2020 20:09
Operator : CG/JU
Sample : L1871-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-17-SA

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.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 |
| 2-Pentanone, 4-hy... | 4.92 | 10.2 | ng | 295262 | 1 | 6.72 | 576561 | 20.0 |
| 5H-Indeno[1,2-b]p... | 11.48 | 2.3 | ng | 121527 | 4 | 11.24 | 1069390 | 20.0 |
| Phenanthrene, 1-m... | 11.74 | 2.3 | ng | 120974 | 4 | 11.24 | 1069390 | 20.0 |
| Anthracene, 2-met... | 11.77 | 6.7 | ng | 359146 | 4 | 11.24 | 1069390 | 20.0 |
| 4H-Cyclopenta[def... | 11.85 | 4.9 | ng | 264061 | 4 | 11.24 | 1069390 | 20.0 |
| Pyrene, 2-methyl- | 13.02 | 2.3 | ng | 169414 | 5 | 13.88 | 1478090 | 20.0 |
| Pyrene, 1-methyl- | 13.12 | 3.0 | ng | 223055 | 5 | 13.88 | 1478090 | 20.0 |
| Cyclotetradecane | 13.72 | 3.5 | ng | 255893 | 5 | 13.88 | 1478090 | 20.0 |
| Benz[a]anthracene... | 14.27 | 2.0 | ng | 150230 | 5 | 13.88 | 1478090 | 20.0 |
| 4,5-Dihydrobenzo[... | 14.76 | 2.4 | ng | 86907 | 6 | 15.31 | 730381 | 20.0 |
| Benzo[e]pyrene | 15.02 | 2.7 | ng | 99531 | 6 | 15.31 | 730381 | 20.0 |
| Perylene | 15.20 | 6.2 | ng | 225550 | 6 | 15.31 | 730381 | 20.0 |
| Heptadecane, 2,6,... | 15.70 | 2.1 | ng | 77393 | 6 | 15.31 | 730381 | 20.0 |
| unknown15.79 | 15.79 | 2.5 | ng | 90888 | 6 | 15.31 | 730381 | 20.0 |