LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

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:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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 | 2.305 | 31 | 37 | 45 | rVB | 109041 | 141329 | 4.02% | 0.397% |
| 2 | 4.116 | 341 | 345 | 354 | rBV | 58815 | 75909 | 2.16% | 0.213% |
| 3 | 5.222 | 529 | 533 | 543 | rBV | 126633 | 168762 | 4.80% | 0.474% |
| 4 | 5.604 | 592 | 598 | 601 | rBV | 3185995 | 3356593 | 95.46% | 9.433% |
| 5 | 6.604 | 762 | 768 | 771 | rBV | 3055984 | 3327965 | 94.64% | 9.352% |
| 6 | 6.745 | 787 | 792 | 796 | rBV | 3252873 | 3483719 | 99.07% | 9.790% |
| 7 | 6.975 | 827 | 831 | 834 | rBV | 1010988 | 783923 | 22.29% | 2.203% |
| 8 | 7.134 | 854 | 858 | 861 | rBV | 2993501 | 2604673 | 74.07% | 7.320% |
| 9 | 7.540 | 922 | 927 | 930 | rBV | 2203144 | 2143483 | 60.96% | 6.023% |
| 10 | 8.257 | 1045 | 1049 | 1051 | rBV | 1280437 | 1019705 | 29.00% | 2.866% |
| 11 | 8.275 | 1051 | 1052 | 1059 | rVB | 195102 | 167238 | 4.76% | 0.470% |
| 12 | 8.969 | 1167 | 1170 | 1177 | rBV | 101784 | 81503 | 2.32% | 0.229% |
| 13 | 9.069 | 1181 | 1187 | 1194 | rBV | 83956 | 78646 | 2.24% | 0.221% |
| 14 | 9.334 | 1227 | 1232 | 1235 | rBV | 3789230 | 3516293 | 100.00% | 9.881% |
| 15 | 9.598 | 1273 | 1277 | 1280 | rBV3 | 32662 | 37874 | 1.08% | 0.106% |
| 16 | 9.669 | 1285 | 1289 | 1292 | rVV | 52974 | 56728 | 1.61% | 0.159% |
| 17 | 9.722 | 1295 | 1298 | 1305 | rVV | 352667 | 356345 | 10.13% | 1.001% |
| 18 | 9.786 | 1305 | 1309 | 1318 | rVB2 | 25754 | 41684 | 1.19% | 0.117% |
| 19 | 9.875 | 1321 | 1324 | 1329 | rBV | 102584 | 93253 | 2.65% | 0.262% |
| 20 | 9.928 | 1329 | 1333 | 1337 | rBV | 92982 | 72155 | 2.05% | 0.203% |
| 21 | 10.016 | 1345 | 1348 | 1352 | rBV | 1292770 | 1085072 | 30.86% | 3.049% |
| 22 | 10.428 | 1415 | 1418 | 1421 | rVB | 105282 | 91766 | 2.61% | 0.258% |
| 23 | 10.563 | 1438 | 1441 | 1448 | rVB | 73831 | 79876 | 2.27% | 0.224% |
| 24 | 10.645 | 1448 | 1455 | 1458 | rBV4 | 33702 | 50374 | 1.43% | 0.142% |
| 25 | 10.810 | 1478 | 1483 | 1486 | rBV | 2279090 | 2196710 | 62.47% | 6.173% |
| 26 | 10.904 | 1495 | 1499 | 1508 | rVB | 93303 | 122543 | 3.49% | 0.344% |
| 27 | 11.110 | 1529 | 1534 | 1537 | rBV3 | 31778 | 50461 | 1.44% | 0.142% |
| 28 | 11.351 | 1572 | 1575 | 1578 | rBV | 67090 | 53008 | 1.51% | 0.149% |
| 29 | 11.404 | 1582 | 1584 | 1587 | rVB | 62178 | 46672 | 1.33% | 0.131% |
| 30 | 11.510 | 1598 | 1602 | 1604 | rBV | 1136417 | 991279 | 28.19% | 2.786% |
| 31 | 11.533 | 1604 | 1606 | 1610 | rVB | 504400 | 365546 | 10.40% | 1.027% |
| 32 | 11.586 | 1612 | 1615 | 1618 | rBV | 109448 | 95068 | 2.70% | 0.267% |
| 33 | 11.839 | 1655 | 1658 | 1661 | rBV | 47913 | 41102 | 1.17% | 0.116% |
| 34 | 12.010 | 1684 | 1687 | 1690 | rBV | 109624 | 122988 | 3.50% | 0.346% |

LSC Area Percent Report

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Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

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:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 12.039 | 1690 | 1692 | 1697 | rVB | 119508 | 96274 | 2.74% | 0.271% |
| 36 | 12.122 | 1703 | 1706 | 1709 | rBV2 | 136020 | 163821 | 4.66% | 0.460% |
| 37 | 12.145 | 1709 | 1710 | 1715 | rVB | 72744 | 49202 | 1.40% | 0.138% |
| 38 | 12.304 | 1734 | 1737 | 1740 | rBV | 62980 | 57856 | 1.65% | 0.163% |
| 39 | 12.339 | 1740 | 1743 | 1749 | rVB2 | 41947 | 43365 | 1.23% | 0.122% |
| 40 | 12.563 | 1777 | 1781 | 1784 | rBV | 85395 | 99574 | 2.83% | 0.280% |
| 41 | 12.651 | 1793 | 1796 | 1799 | rBV3 | 34549 | 45669 | 1.30% | 0.128% |
| 42 | 12.727 | 1805 | 1809 | 1813 | rBV | 529603 | 432461 | 12.30% | 1.215% |
| 43 | 12.822 | 1822 | 1825 | 1828 | rBV | 81676 | 67920 | 1.93% | 0.191% |
| 44 | 12.898 | 1836 | 1838 | 1841 | rVB2 | 63640 | 51020 | 1.45% | 0.143% |
| 45 | 12.957 | 1845 | 1848 | 1854 | rVB | 484130 | 479606 | 13.64% | 1.348% |
| 46 | 13.098 | 1867 | 1872 | 1875 | rBV | 2920135 | 2676594 | 76.12% | 7.522% |
| 47 | 13.169 | 1881 | 1884 | 1891 | rBV2 | 43839 | 71696 | 2.04% | 0.201% |
| 48 | 13.263 | 1897 | 1900 | 1901 | rBV3 | 43360 | 45329 | 1.29% | 0.127% |
| 49 | 13.286 | 1901 | 1904 | 1910 | rVB2 | 118456 | 121224 | 3.45% | 0.341% |
| 50 | 13.351 | 1913 | 1915 | 1918 | rVB2 | 78749 | 59604 | 1.70% | 0.167% |
| 51 | 13.392 | 1918 | 1922 | 1926 | rBV2 | 70294 | 71147 | 2.02% | 0.200% |
| 52 | 13.516 | 1940 | 1943 | 1946 | rBV | 50379 | 44344 | 1.26% | 0.125% |
| 53 | 13.639 | 1961 | 1964 | 1968 | rBV4 | 23122 | 45294 | 1.29% | 0.127% |
| 54 | 13.910 | 2005 | 2010 | 2012 | rBV2 | 60891 | 74572 | 2.12% | 0.210% |
| 55 | 13.974 | 2017 | 2021 | 2034 | rVB4 | 505764 | 617934 | 17.57% | 1.736% |
| 56 | 14.163 | 2048 | 2053 | 2056 | rBV2 | 925750 | 1054500 | 29.99% | 2.963% |
| 57 | 14.186 | 2056 | 2057 | 2065 | rVB | 215149 | 165127 | 4.70% | 0.464% |
| 58 | 14.557 | 2115 | 2120 | 2123 | rBV | 50022 | 61335 | 1.74% | 0.172% |
| 59 | 14.621 | 2128 | 2131 | 2135 | rBV4 | 46367 | 64640 | 1.84% | 0.182% |
| 60 | 15.245 | 2233 | 2237 | 2241 | rBV | 138755 | 244305 | 6.95% | 0.687% |
| 61 | 15.292 | 2243 | 2245 | 2249 | rVB2 | 130138 | 146410 | 4.16% | 0.411% |
| 62 | 15.363 | 2254 | 2257 | 2265 | rVB | 49124 | 82201 | 2.34% | 0.231% |
| 63 | 15.574 | 2289 | 2293 | 2299 | rVB | 98605 | 144257 | 4.10% | 0.405% |
| 64 | 15.639 | 2300 | 2304 | 2310 | rVB | 136766 | 175082 | 4.98% | 0.492% |
| 65 | 15.710 | 2310 | 2316 | 2320 | rBV | 544403 | 731843 | 20.81% | 2.057% |
| 66 | 16.168 | 2390 | 2394 | 2399 | rVB3 | 45703 | 63436 | 1.80% | 0.178% |
| 67 | 16.221 | 2400 | 2403 | 2410 | rVV4 | 45993 | 79978 | 2.27% | 0.225% |
| 68 | 17.286 | 2579 | 2584 | 2590 | rBV2 | 50917 | 104238 | 2.96% | 0.293% |
| 69 | 17.768 | 2662 | 2666 | 2670 | rVB | 31280 | 53245 | 1.51% | 0.150% |

Sum of corrected areas: 35585348

LSC Report - Integrated Chromatogram

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Data File : BF103948.D
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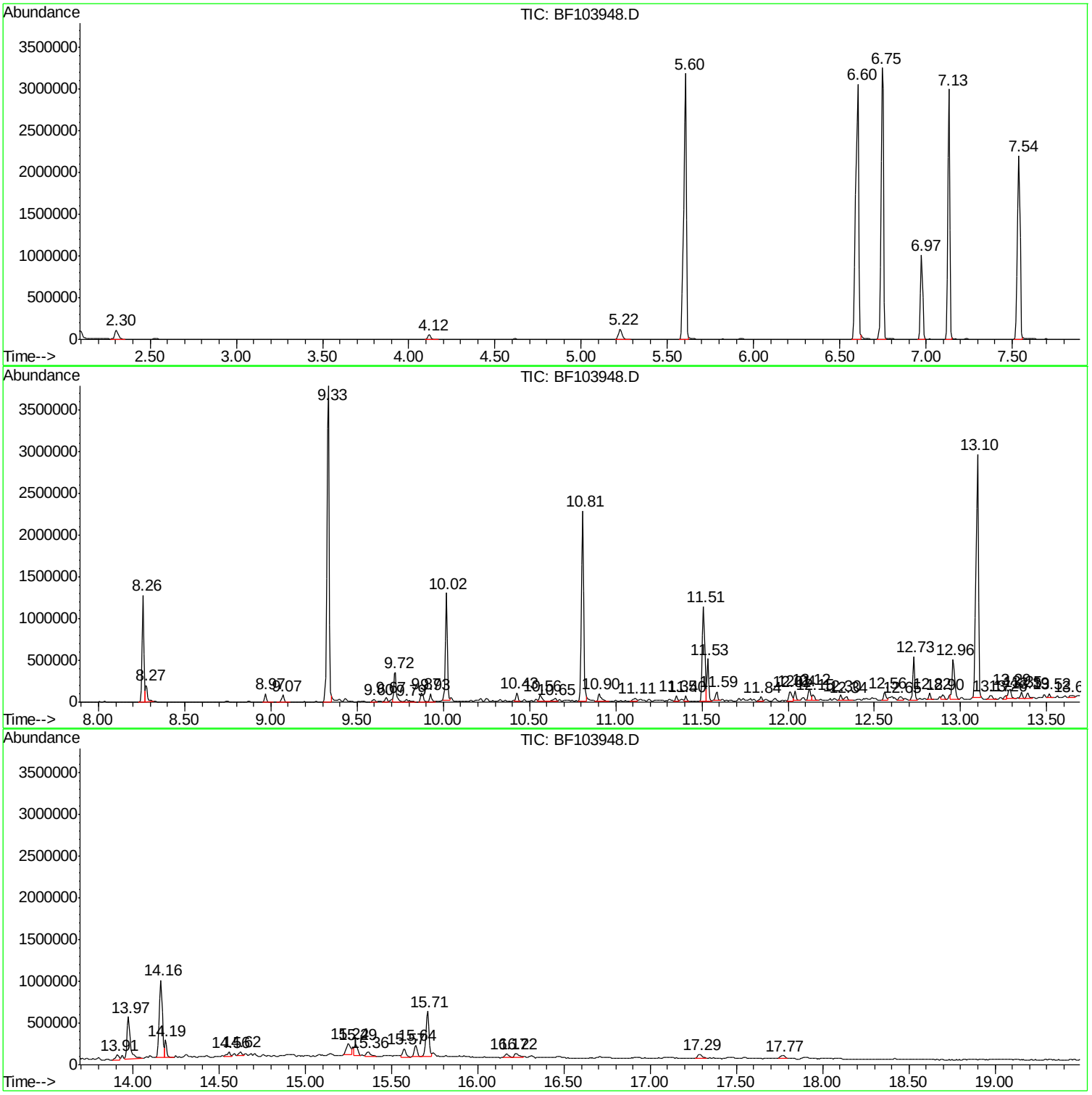
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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**Instrument :** BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
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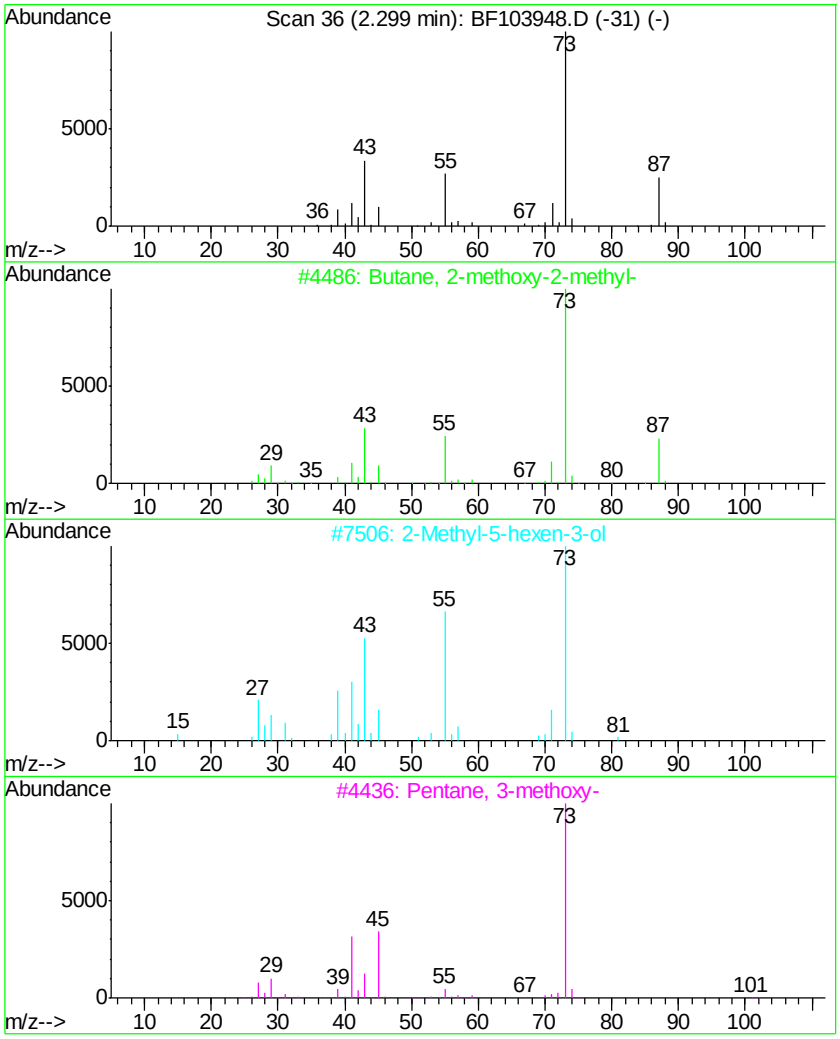
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

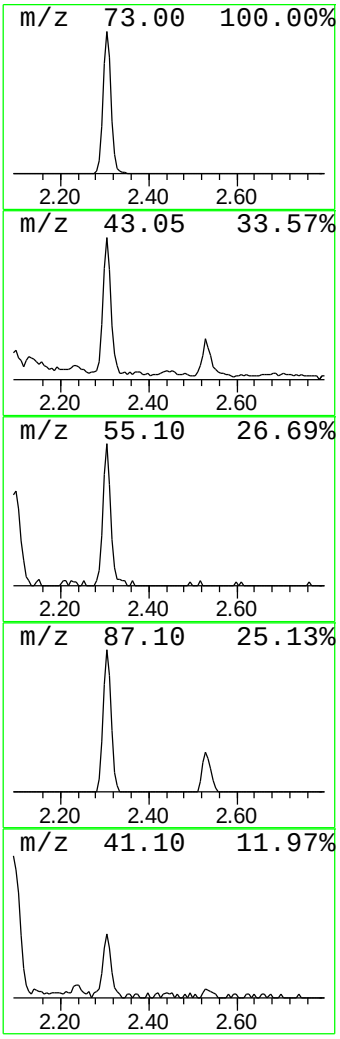
**********

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.30 | 3.61 ng | 141329 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 64 |
| 2 | 2-Methyl-5-hexen-3-ol | 114 | C7H14O | 032815-70-6 | 38 |
| 3 | Pentane, 3-methoxy- | 102 | C6H14O | 036839-67-5 | 17 |
| 4 | 3-Hexanol, 2-methyl- | 116 | C7H16O | 000617-29-8 | 17 |
| 5 | N-Ethylformamide | 73 | C3H7NO | 000627-45-2 | 9 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

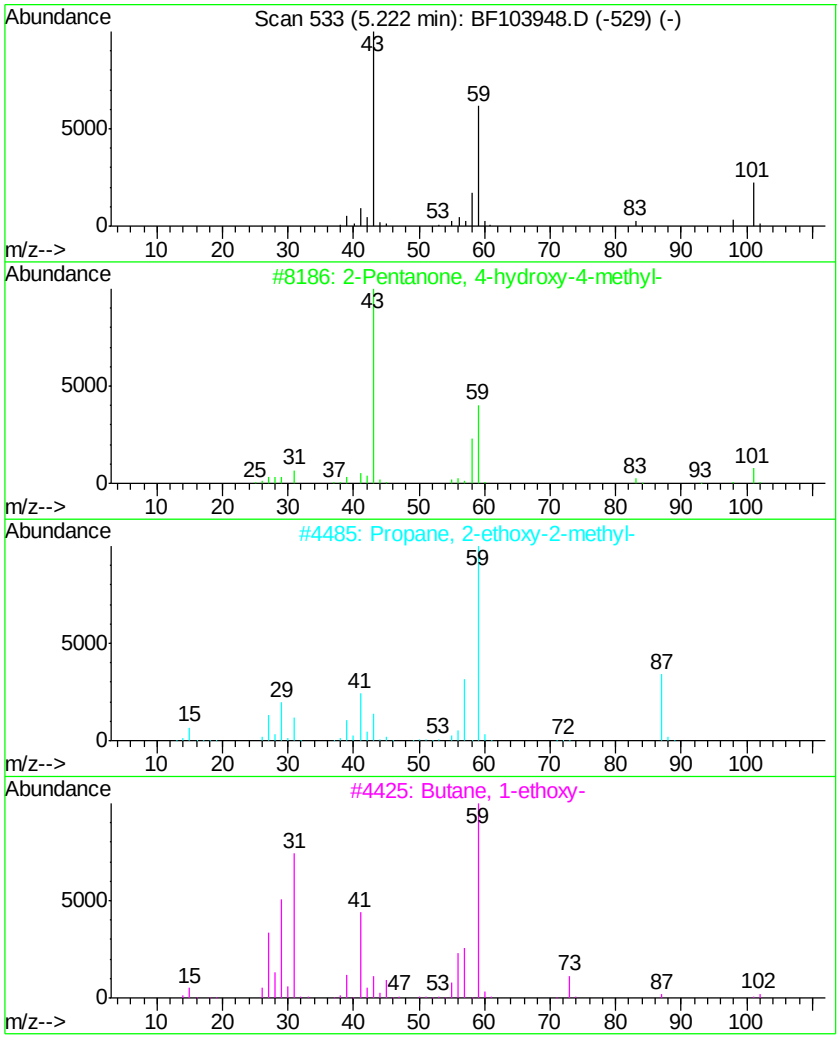
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

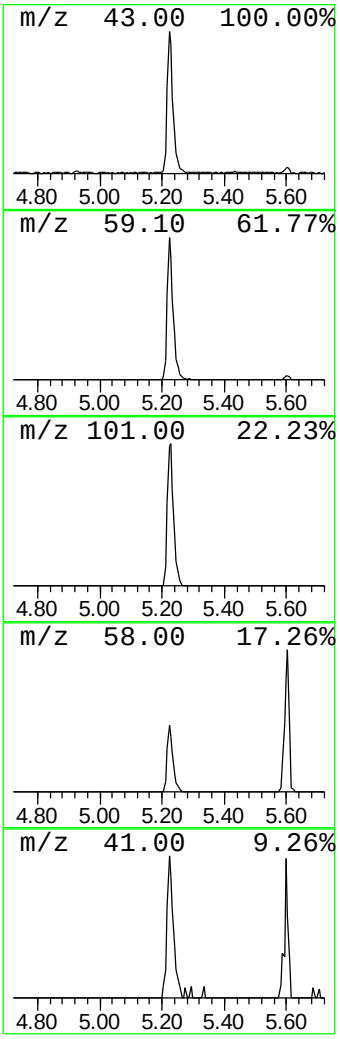
**********

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

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.22 | 4.31 ng | 168762 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 56 |
| 2 | Propane, 2-ethoxy-2-methyl- | 102 | C6H14O | 000637-92-3 | 17 |
| 3 | Butane, 1-ethoxy- | 102 | C6H14O | 000628-81-9 | 9 |
| 4 | Morpholine, 4-methyl- | 101 | C5H11NO | 000109-02-4 | 9 |
| 5 | Propanamide, N-ethyl- | 101 | C5H11NO | 005129-72-6 | 9 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

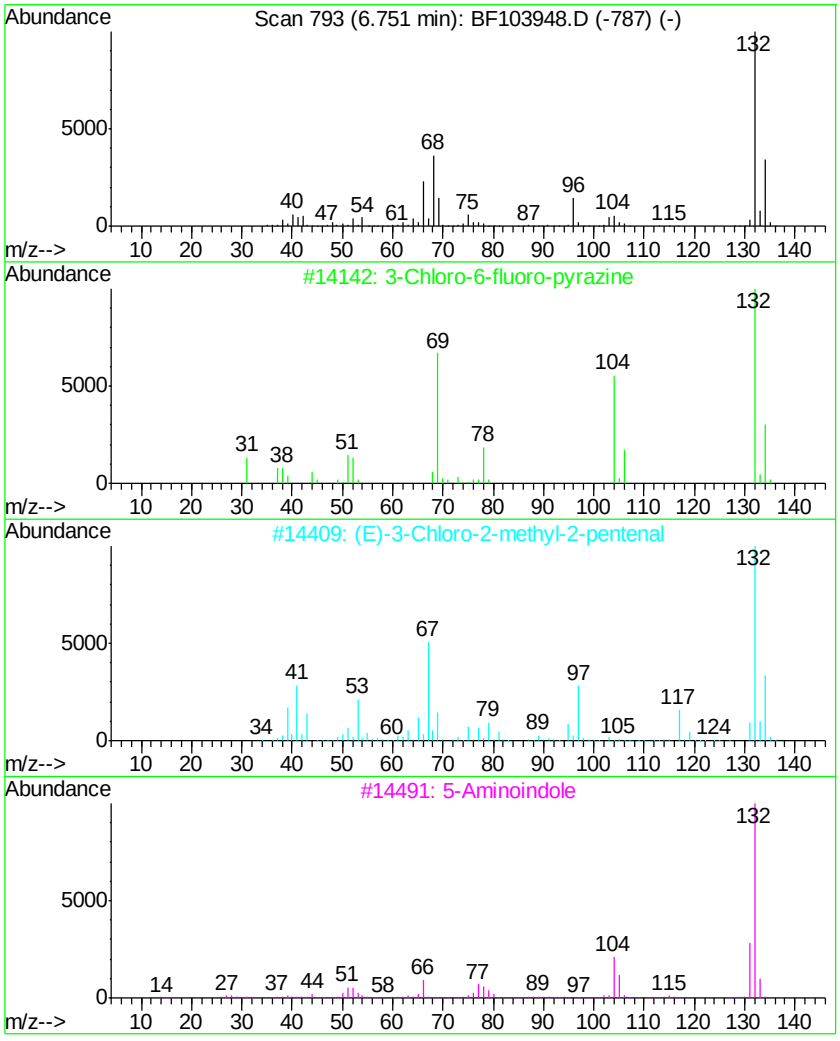
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

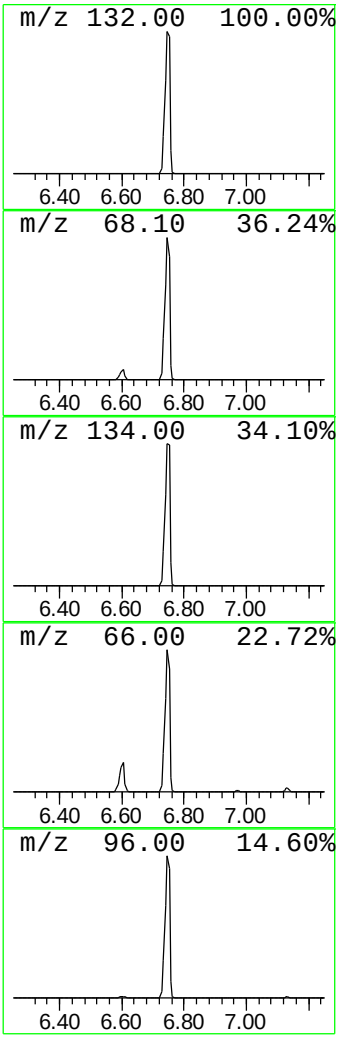
**********

Peak Number 3 unknown6.75 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.75 | 88.88 ng | 3483720 | 1,4-Dichlorobenzene-d4 | 6.97 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Chloro-6-fluoro-pyrazine | 132 | C4H2ClFN2 | 1000146-10-7 | 27 |
| 2 | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO | 031357-76-3 | 17 |
| 3 | 5-Aminoindole | 132 | C8H8N2 | 005192-03-0 | 12 |
| 4 | Tranylcypromine-propionyl | 189 | C12H15NO | 1000123-86-3 | 10 |
| 5 | 4-Chloro-2-fluoro-pyrimidine | 132 | C4H2ClFN2 | 051422-00-5 | 9 |



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**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

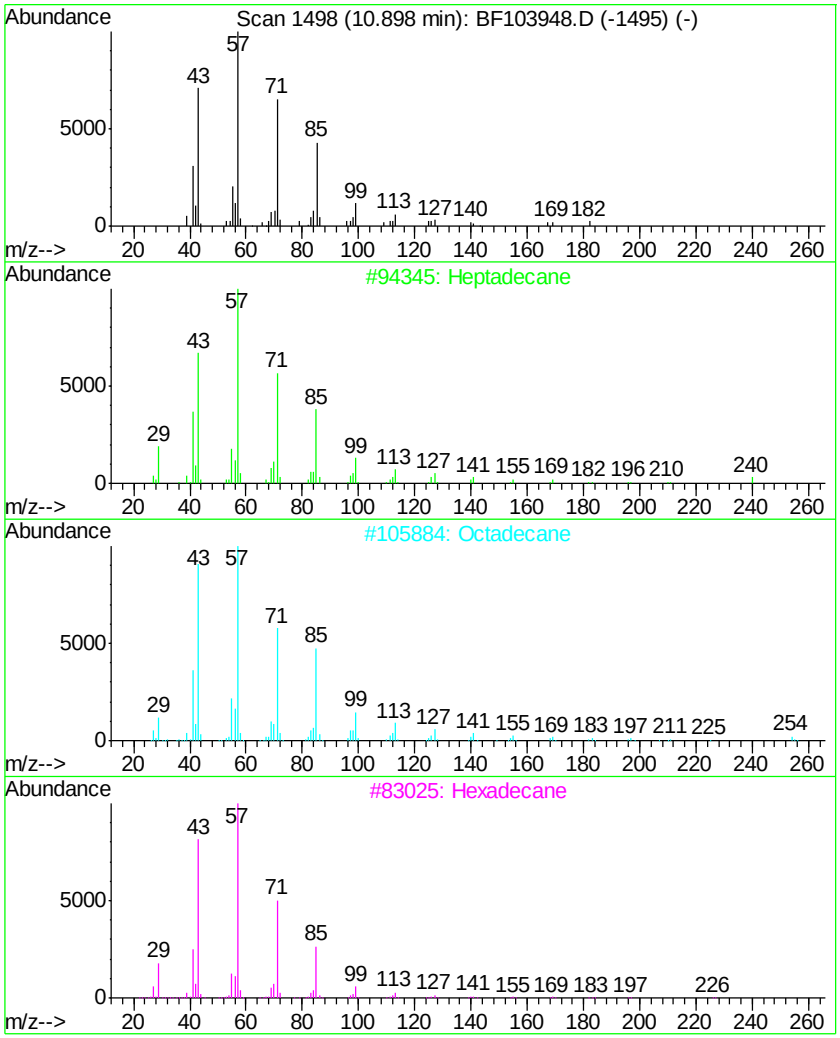
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

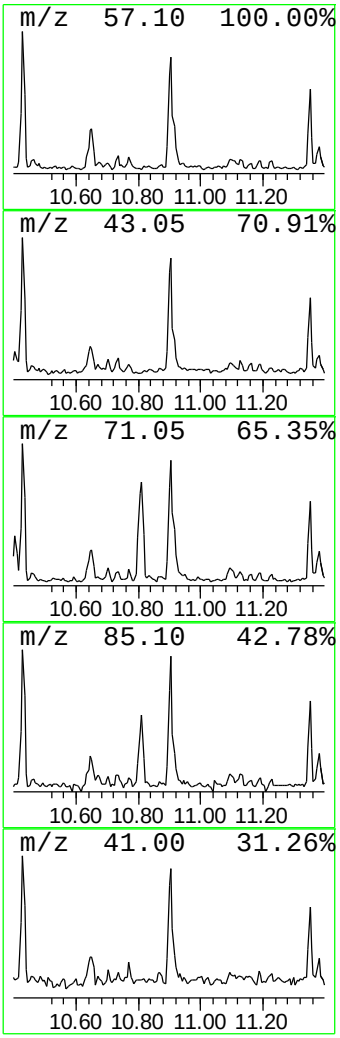
**********

Peak Number 5 Heptadecane Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.90 | 2.47 ng | 122543 | Phenanthrene-d10 | 11.51 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptadecane | 240 | C17H36 | 000629-78-7 | 91 |
| 2 | Octadecane | 254 | C18H38 | 000593-45-3 | 91 |
| 3 | Hexadecane | 226 | C16H34 | 000544-76-3 | 90 |
| 4 | Heneicosane | 296 | C21H44 | 000629-94-7 | 87 |
| 5 | Tetradecane | 198 | C14H30 | 000629-59-4 | 86 |



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Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

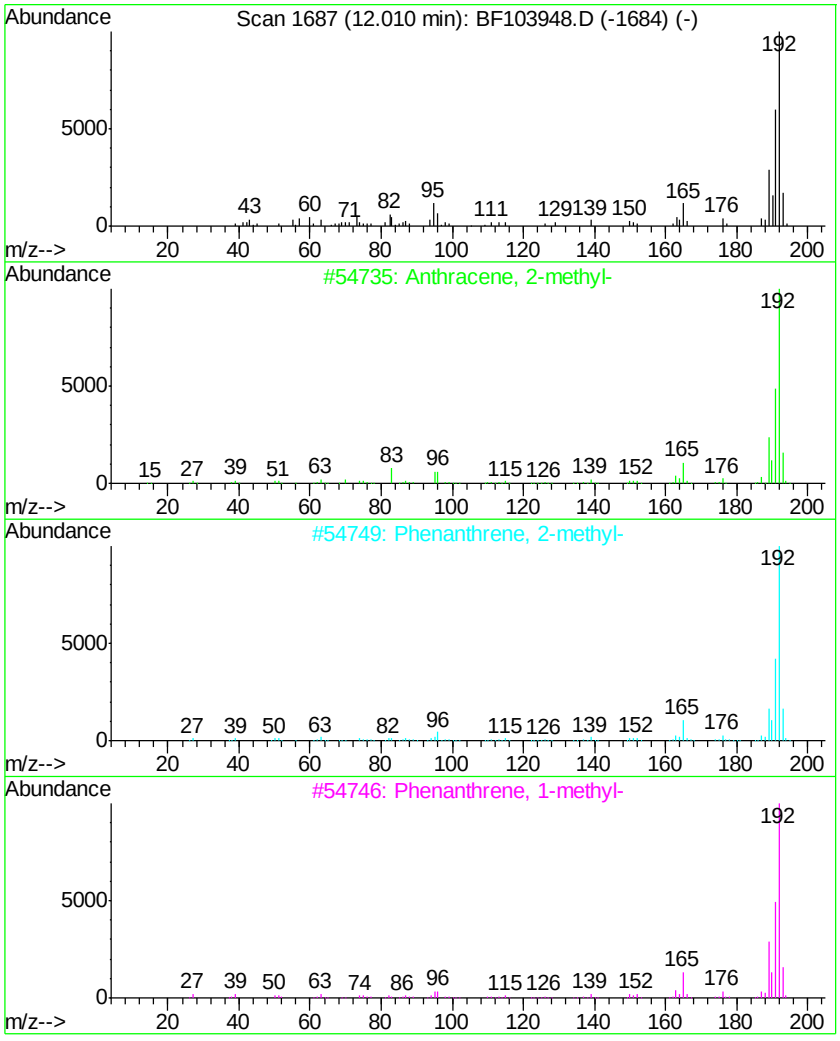
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

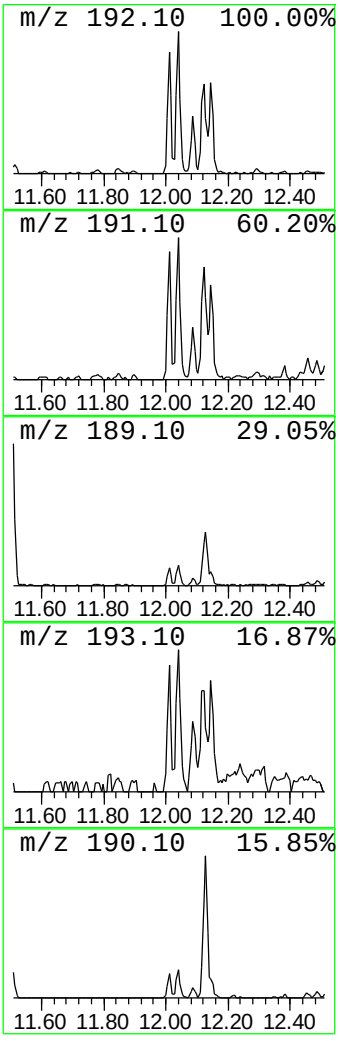
**********

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.01 | 2.48 ng | 122988 | Phenanthrene-d10 | 11.51 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 96 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 95 |
| 3 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 94 |
| 4 | 1H-Indene, 2-phenyl- | 192 | C15H12 | 004505-48-0 | 93 |
| 5 | 5H-Dibenzo[a,d]cycloheptene | 192 | C15H12 | 000256-81-5 | 93 |



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Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

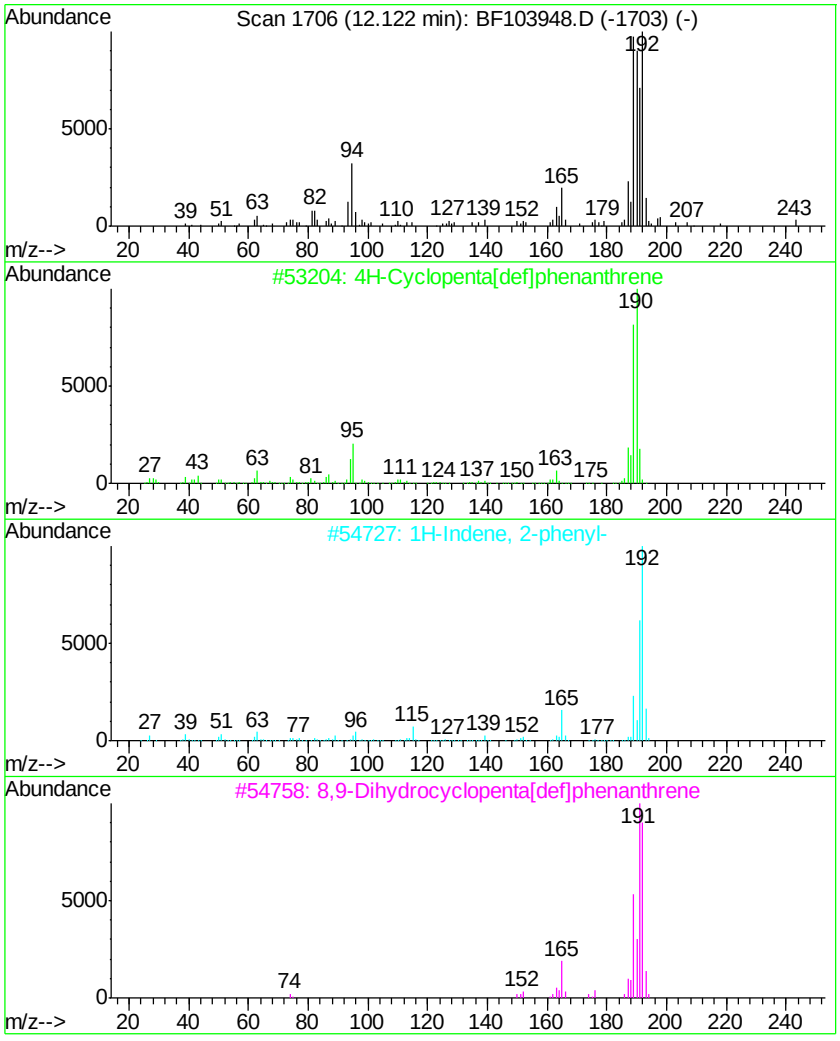
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

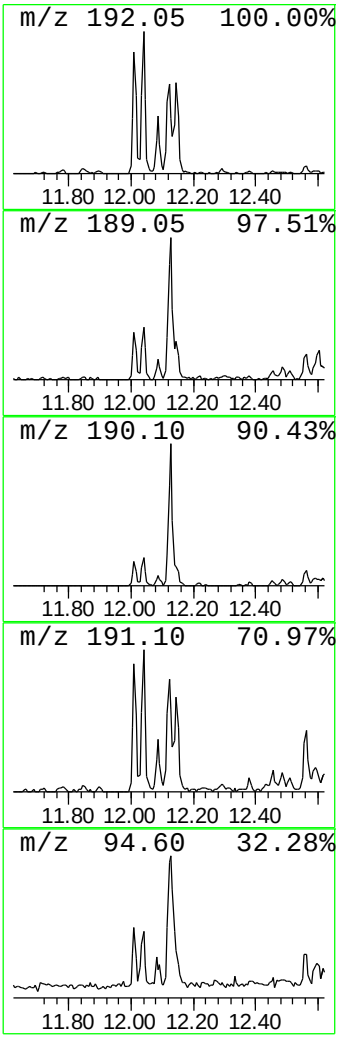
**********

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.12 | 3.31 ng | 163821 | Phenanthrene-d10 | 11.51 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 50 |
| 2 | 1H-Indene, 2-phenyl- | 192 | C15H12 | 004505-48-0 | 46 |
| 3 | 8,9-Dihydrocyclopenta[def]phenan... | 192 | C15H12 | 027410-55-5 | 44 |
| 4 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 42 |
| 5 | Phenanthrene, 4-methyl- | 192 | C15H12 | 000832-64-4 | 42 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

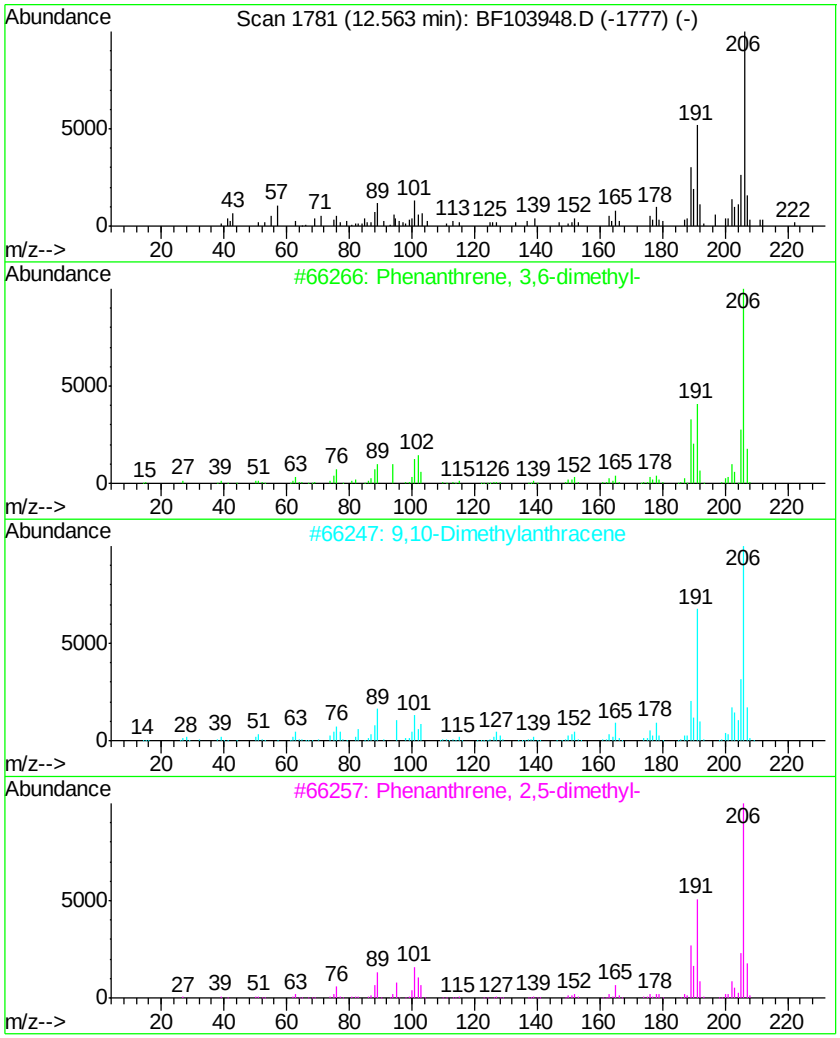
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

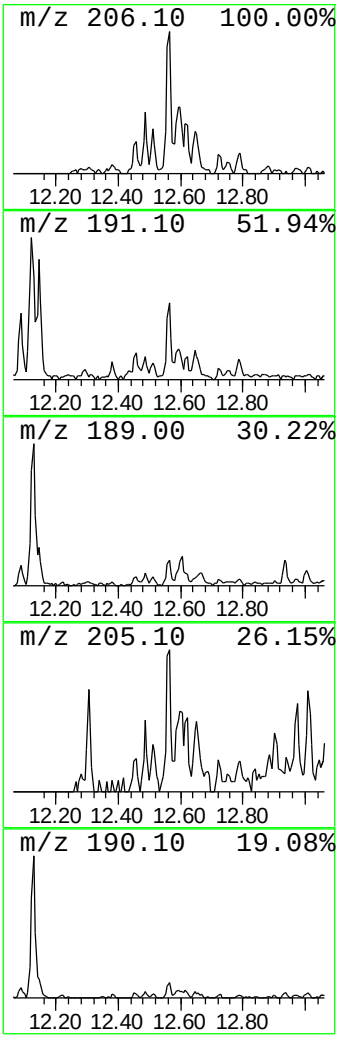
**********

Peak Number 8 Phenanthrene, 3,6-dimethyl- Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.56 | 2.01 ng | 99574 | Phenanthrene-d10 | 11.51 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 93 |
| 2 | 9,10-Dimethylanthracene | 206 | C16H14 | 000781-43-1 | 93 |
| 3 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 93 |
| 4 | Phenanthrene, 2,7-dimethyl- | 206 | C16H14 | 001576-69-8 | 89 |
| 5 | Anthracene, 1,4-dimethyl- | 206 | C16H14 | 000781-92-0 | 87 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

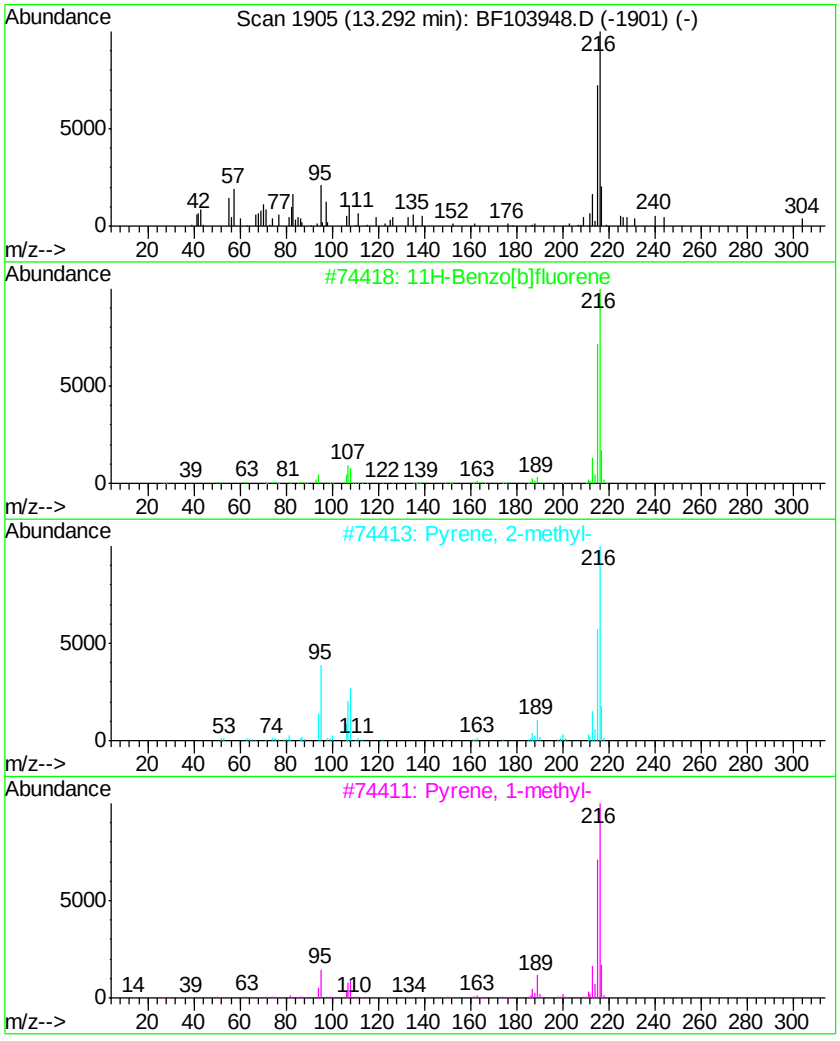
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

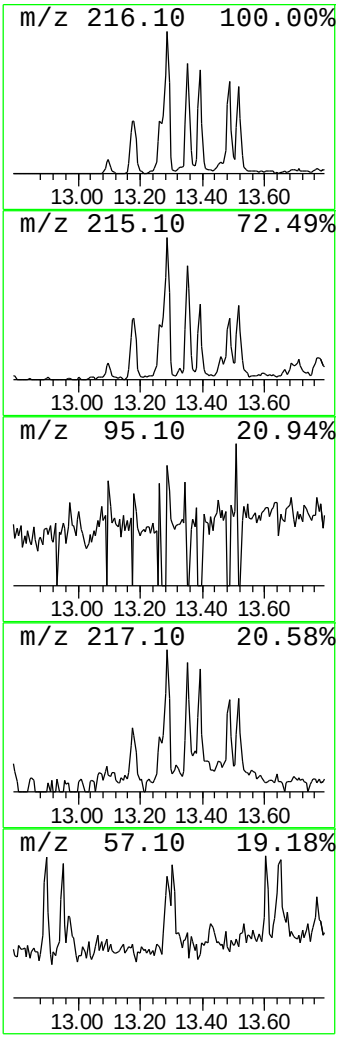
**********

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.29 | 2.30 ng | 121224 | Chrysene-d12 | 14.16 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 93 |
| 2 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 91 |
| 3 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 91 |
| 4 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 87 |
| 5 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 76 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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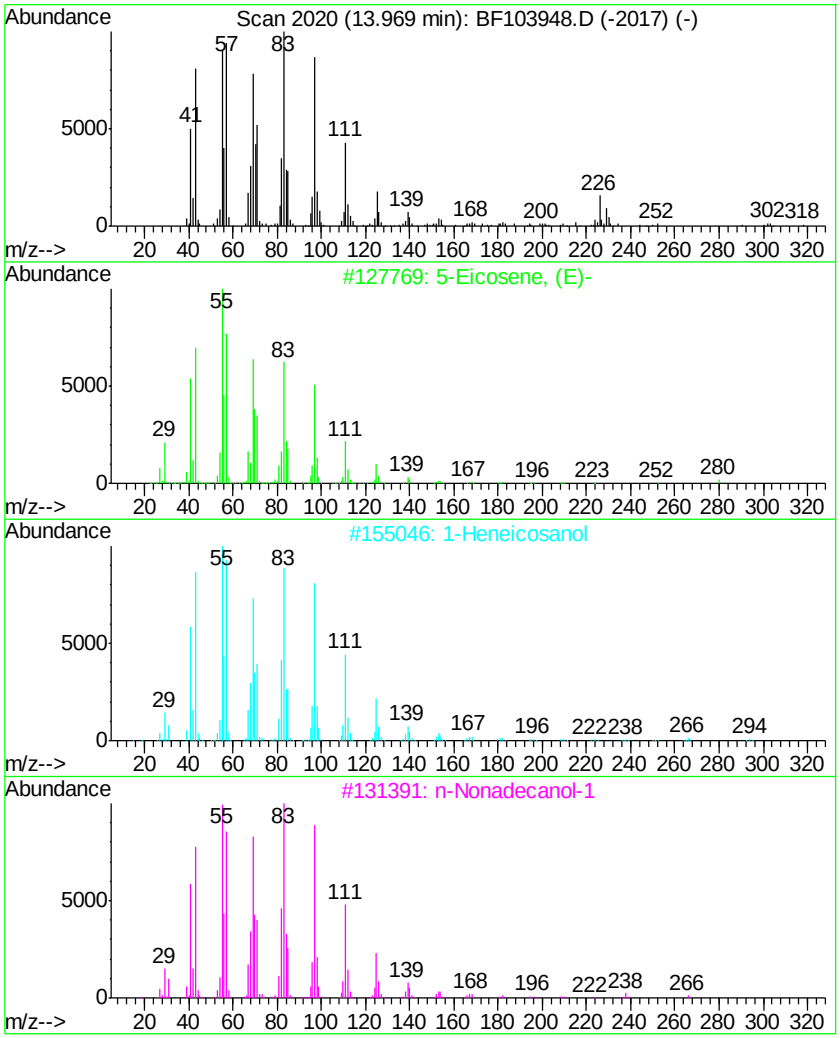
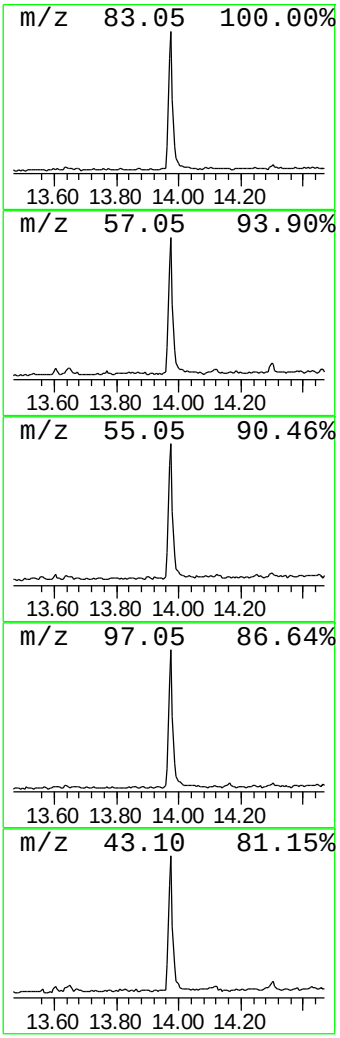
Peak Number 10 5-Eicosene, (E)- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.97 | 11.72 ng | 617934 | Chrysene-d12 | 14.16 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 5-Eicosene, (E)- | 280 | C20H40 | 074685-30-6 | 99 |
| 2 | 1-Heneicosanol | 312 | C21H44O | 015594-90-8 | 95 |
| 3 | n-Nonadecanol-1 | 284 | C19H40O | 001454-84-8 | 95 |
| 4 | Trifluoroacetic acid, pentadecyl... | 324 | C17H31F3O2 | 959010-23-2 | 94 |
| 5 | Heptafluorobutyric acid, pentade... | 424 | C19H31F7O2 | 959261-23-5 | 94 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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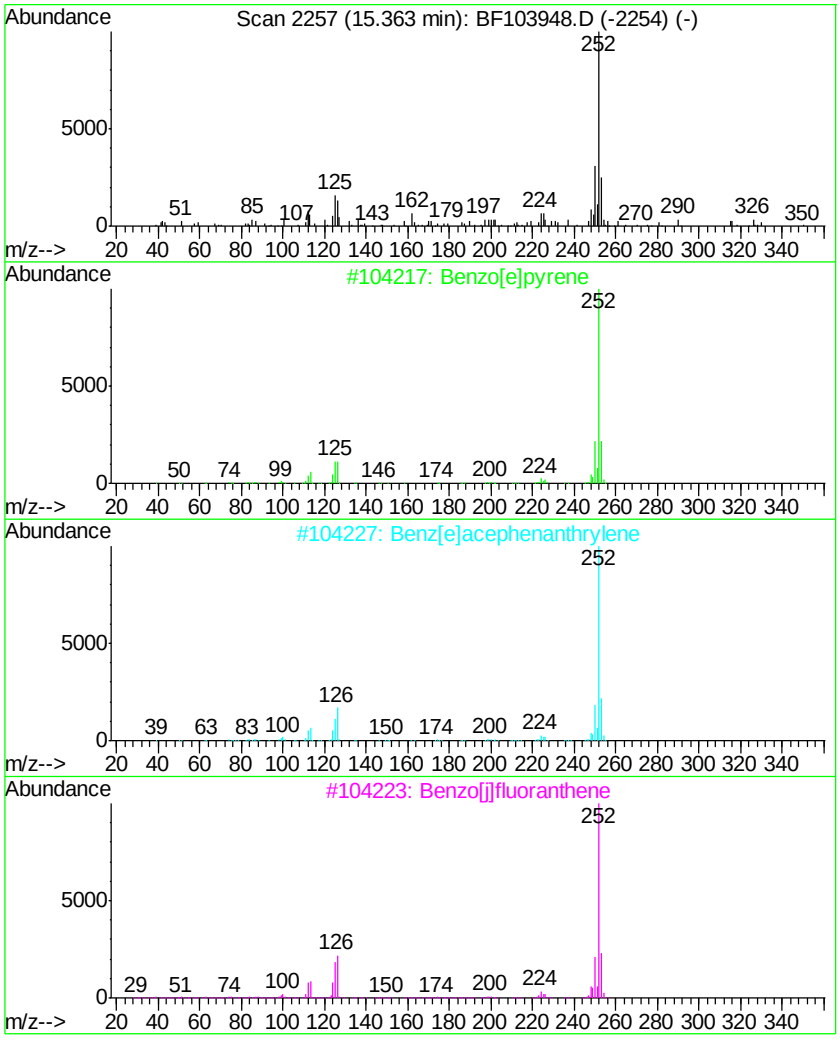
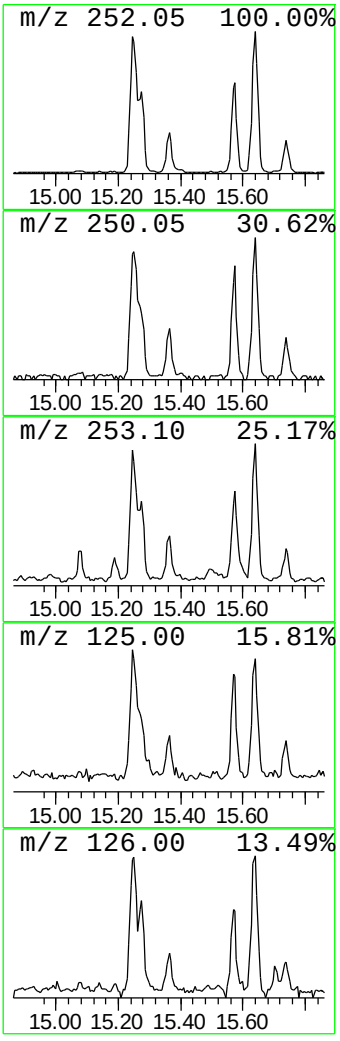
Peak Number 11 Benzo[e]pyrene Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.36 | 2.25 ng | 82201 | Perylene-d12 | 15.71 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 95 |
| 2 | Benz[e]acephenanthrylene | 252 | C20H12 | 000205-99-2 | 89 |
| 3 | Benzo[j]fluoranthene | 252 | C20H12 | 000205-82-3 | 76 |
| 4 | Benzo[a]pyrene | 252 | C20H12 | 000050-32-8 | 76 |
| 5 | Perylene | 252 | C20H12 | 000198-55-0 | 76 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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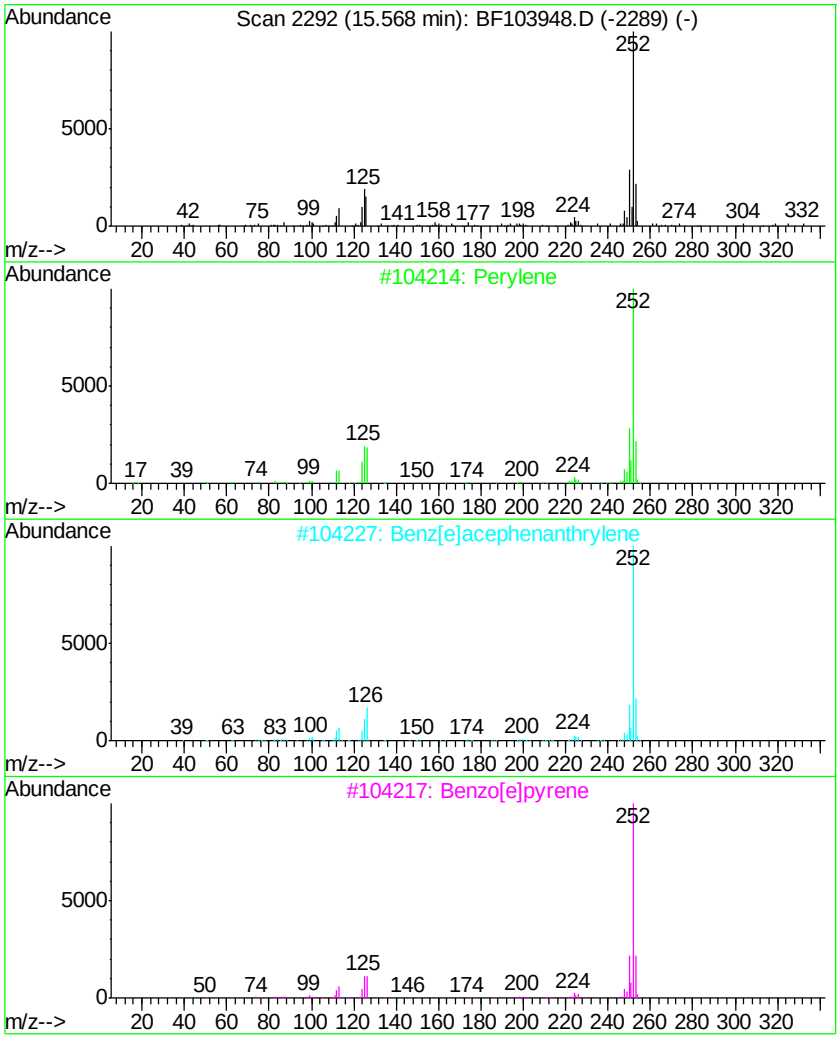
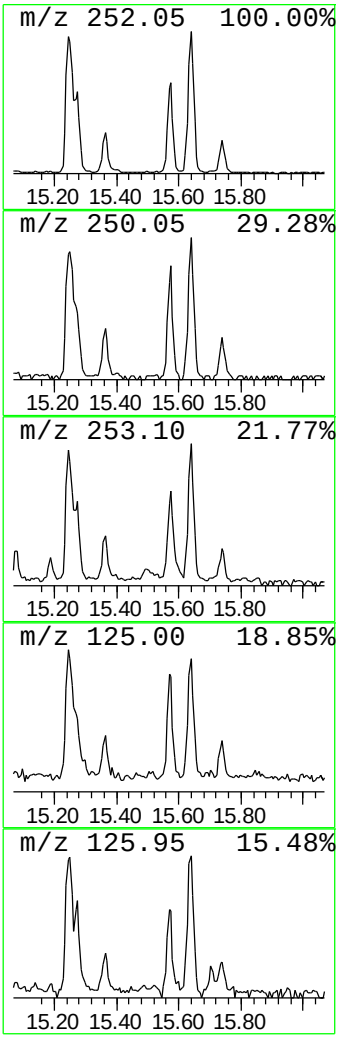
Peak Number 12 Perylene Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.57 | 3.94 ng | 144257 | Perylene-d12 | 15.71 |

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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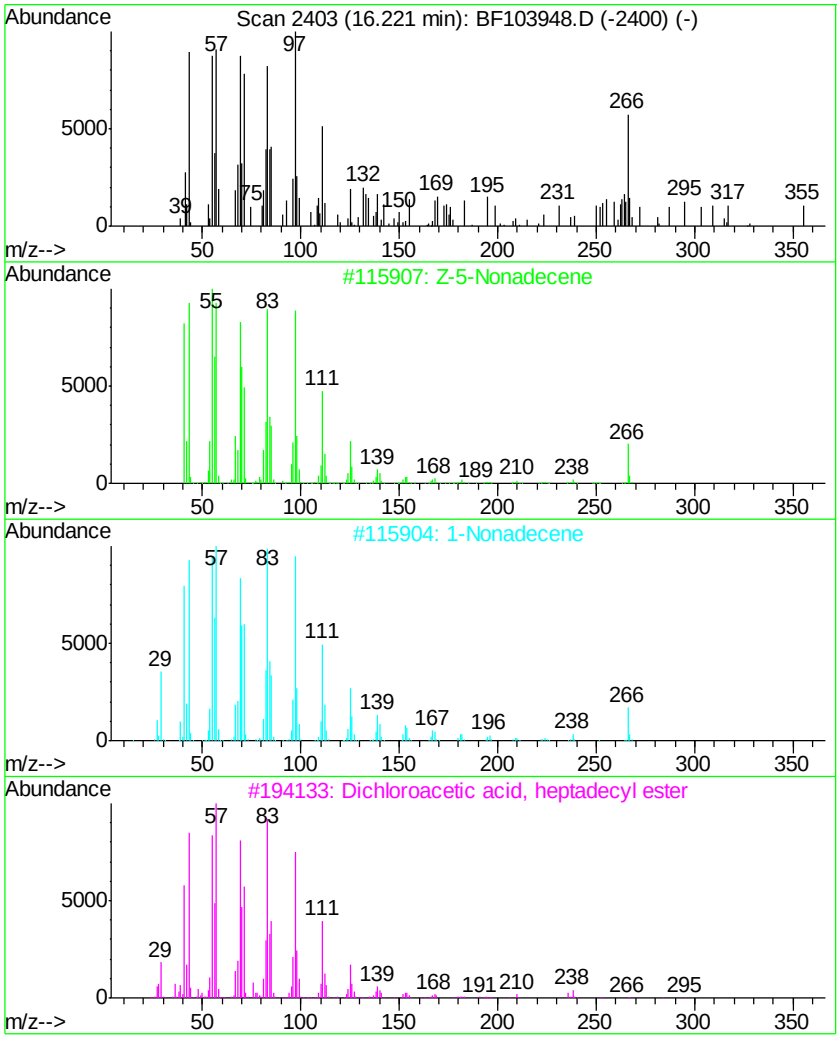
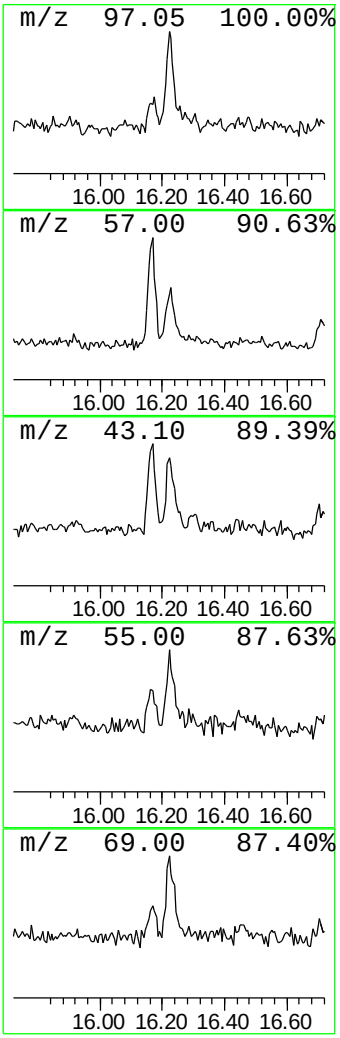
Peak Number 13 Z-5-Nonadecene Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.22 | 2.19 ng | 79978 | Perylene-d12 | 15.71 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Z-5-Nonadecene | 266 | C19H38 | 1000131-11-8 | 83 |
| 2 | 1-Nonadecene | 266 | C19H38 | 018435-45-5 | 83 |
| 3 | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 74 |
| 4 | 1-Heptacosanol | 396 | C27H56O | 002004-39-9 | 68 |
| 5 | Tetradecyl trifluoroacetate | 310 | C16H29F3O2 | 006222-02-2 | 64 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103948.D
Acq On : 27 Mar 2018 1:35
Operator : JU/SJ
Sample : J1998-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** SB-101(9-10)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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... | 2.30 | 3.6 | ng | 141329 | 1 | 6.97 | 783923 | 20.0 |
| 2-Pentanone, 4-hy... | 5.22 | 4.3 | ng | 168762 | 1 | 6.97 | 783923 | 20.0 |
| unknown6.75 | 6.75 | 88.9 | ng | 3483720 | 1 | 6.97 | 783923 | 20.0 |
| Heptadecane | 10.90 | 2.5 | ng | 122543 | 4 | 11.51 | 991279 | 20.0 |
| Anthracene, 2-met... | 12.01 | 2.5 | ng | 122988 | 4 | 11.51 | 991279 | 20.0 |
| 4H-Cyclopenta[def... | 12.12 | 3.3 | ng | 163821 | 4 | 11.51 | 991279 | 20.0 |
| Phenanthrene, 3,6... | 12.56 | 2.0 | ng | 99574 | 4 | 11.51 | 991279 | 20.0 |
| 11H-Benzo[b]fluorene | 13.29 | 2.3 | ng | 121224 | 5 | 14.16 | 1054500 | 20.0 |
| 5-Eicosene, (E)- | 13.97 | 11.7 | ng | 617934 | 5 | 14.16 | 1054500 | 20.0 |
| Benzo[e]pyrene | 15.36 | 2.3 | ng | 82201 | 6 | 15.71 | 731843 | 20.0 |
| Perylene | 15.57 | 3.9 | ng | 144257 | 6 | 15.71 | 731843 | 20.0 |
| Z-5-Nonadecene | 16.22 | 2.2 | ng | 79978 | 6 | 15.71 | 731843 | 20.0 |