LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

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-BF040618.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 | 3.834 | 291 | 297 | 304 | rBV | 167546 | 234793 | 8.79% | 0.930% |
| 2 | 5.022 | 493 | 499 | 504 | rBV | 97970 | 141669 | 5.30% | 0.561% |
| 3 | 5.434 | 565 | 569 | 572 | rBV | 2305048 | 2144603 | 80.24% | 8.497% |
| 4 | 6.457 | 738 | 743 | 746 | rBV | 2197004 | 2229155 | 83.41% | 8.832% |
| 5 | 6.586 | 760 | 765 | 768 | rBV | 2473925 | 2389102 | 89.39% | 9.466% |
| 6 | 6.810 | 799 | 803 | 806 | rBV | 743523 | 617182 | 23.09% | 2.445% |
| 7 | 6.963 | 825 | 829 | 833 | rBV | 2160460 | 1986513 | 74.33% | 7.871% |
| 8 | 7.375 | 894 | 899 | 902 | rBV | 1492729 | 1449647 | 54.24% | 5.744% |
| 9 | 8.092 | 1018 | 1021 | 1024 | rBV | 972153 | 783295 | 29.31% | 3.104% |
| 10 | 9.169 | 1200 | 1204 | 1207 | rBV | 3099186 | 2672612 | 100.00% | 10.589% |
| 11 | 9.563 | 1267 | 1271 | 1278 | rVB | 213194 | 213441 | 7.99% | 0.846% |
| 12 | 9.710 | 1293 | 1296 | 1303 | rBV | 56856 | 54702 | 2.05% | 0.217% |
| 13 | 9.769 | 1303 | 1306 | 1309 | rVV | 81605 | 62737 | 2.35% | 0.249% |
| 14 | 9.851 | 1316 | 1320 | 1324 | rBV | 984837 | 836053 | 31.28% | 3.313% |
| 15 | 10.269 | 1388 | 1391 | 1394 | rVB | 107318 | 87756 | 3.28% | 0.348% |
| 16 | 10.486 | 1424 | 1428 | 1431 | rBV | 27503 | 32419 | 1.21% | 0.128% |
| 17 | 10.645 | 1451 | 1455 | 1464 | rVB | 1501252 | 1408400 | 52.70% | 5.580% |
| 18 | 10.745 | 1469 | 1472 | 1480 | rVB2 | 88218 | 123543 | 4.62% | 0.489% |
| 19 | 11.192 | 1545 | 1548 | 1551 | rBV3 | 51391 | 43883 | 1.64% | 0.174% |
| 20 | 11.339 | 1569 | 1573 | 1575 | rBV | 868488 | 737990 | 27.61% | 2.924% |
| 21 | 11.363 | 1575 | 1577 | 1583 | rVV | 318191 | 266742 | 9.98% | 1.057% |
| 22 | 11.416 | 1583 | 1586 | 1589 | rVV | 75837 | 62285 | 2.33% | 0.247% |
| 23 | 11.574 | 1609 | 1613 | 1618 | rBV2 | 24003 | 33891 | 1.27% | 0.134% |
| 24 | 11.621 | 1618 | 1621 | 1627 | rVB5 | 20917 | 30524 | 1.14% | 0.121% |
| 25 | 11.739 | 1639 | 1641 | 1647 | rVB2 | 28478 | 34264 | 1.28% | 0.136% |
| 26 | 11.839 | 1656 | 1658 | 1660 | rBV | 56323 | 53087 | 1.99% | 0.210% |
| 27 | 11.868 | 1660 | 1663 | 1666 | rVB2 | 88823 | 102100 | 3.82% | 0.405% |
| 28 | 11.951 | 1674 | 1677 | 1680 | rBV2 | 70788 | 80515 | 3.01% | 0.319% |
| 29 | 11.974 | 1680 | 1681 | 1686 | rVB | 42652 | 36942 | 1.38% | 0.146% |
| 30 | 12.027 | 1686 | 1690 | 1693 | rBV3 | 33130 | 36513 | 1.37% | 0.145% |
| 31 | 12.139 | 1706 | 1709 | 1712 | rBV | 39986 | 35230 | 1.32% | 0.140% |
| 32 | 12.168 | 1712 | 1714 | 1719 | rVB3 | 33885 | 36606 | 1.37% | 0.145% |
| 33 | 12.392 | 1748 | 1752 | 1754 | rBV | 59578 | 62762 | 2.35% | 0.249% |
| 34 | 12.480 | 1764 | 1767 | 1773 | rBV2 | 37225 | 44277 | 1.66% | 0.175% |

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

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-BF040618.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 12.557 | 1777 | 1780 | 1784 | rVB | 547252 | 454749 | 17.02% | 1.802% |
| 36 | 12.786 | 1813 | 1819 | 1825 | rBV | 496038 | 530888 | 19.86% | 2.103% |
| 37 | 12.839 | 1825 | 1828 | 1831 | rVB3 | 29500 | 36477 | 1.36% | 0.145% |
| 38 | 12.927 | 1836 | 1843 | 1847 | rBV | 2190461 | 2142072 | 80.15% | 8.487% |
| 39 | 13.004 | 1852 | 1856 | 1860 | rBV4 | 38854 | 65042 | 2.43% | 0.258% |
| 40 | 13.092 | 1868 | 1871 | 1872 | rBV2 | 41306 | 39201 | 1.47% | 0.155% |
| 41 | 13.115 | 1872 | 1875 | 1878 | rVB | 70962 | 75938 | 2.84% | 0.301% |
| 42 | 13.180 | 1884 | 1886 | 1889 | rBV | 38270 | 33458 | 1.25% | 0.133% |
| 43 | 13.221 | 1889 | 1893 | 1895 | rVV2 | 54212 | 60655 | 2.27% | 0.240% |
| 44 | 13.398 | 1921 | 1923 | 1926 | rVB2 | 42178 | 30377 | 1.14% | 0.120% |
| 45 | 13.598 | 1954 | 1957 | 1959 | rBV3 | 34630 | 44019 | 1.65% | 0.174% |
| 46 | 13.627 | 1959 | 1962 | 1965 | rVB | 61700 | 67630 | 2.53% | 0.268% |
| 47 | 13.733 | 1978 | 1980 | 1983 | rBV2 | 45729 | 42658 | 1.60% | 0.169% |
| 48 | 13.786 | 1987 | 1989 | 1991 | rVV2 | 35339 | 38604 | 1.44% | 0.153% |
| 49 | 13.815 | 1991 | 1994 | 2001 | rVB3 | 125667 | 179532 | 6.72% | 0.711% |
| 50 | 13.957 | 2015 | 2018 | 2019 | rBV | 72638 | 70646 | 2.64% | 0.280% |
| 51 | 13.992 | 2019 | 2024 | 2026 | rVV2 | 725972 | 941064 | 35.21% | 3.729% |
| 52 | 14.015 | 2026 | 2028 | 2036 | rVB | 287744 | 257577 | 9.64% | 1.021% |
| 53 | 15.033 | 2198 | 2201 | 2204 | rBV | 178511 | 243164 | 9.10% | 0.963% |
| 54 | 15.339 | 2250 | 2253 | 2257 | rVB | 105649 | 125625 | 4.70% | 0.498% |
| 55 | 15.398 | 2260 | 2263 | 2269 | rVB | 125934 | 170234 | 6.37% | 0.674% |
| 56 | 15.462 | 2270 | 2274 | 2278 | rBV | 342346 | 423919 | 15.86% | 1.680% |

Sum of corrected areas: 25238762

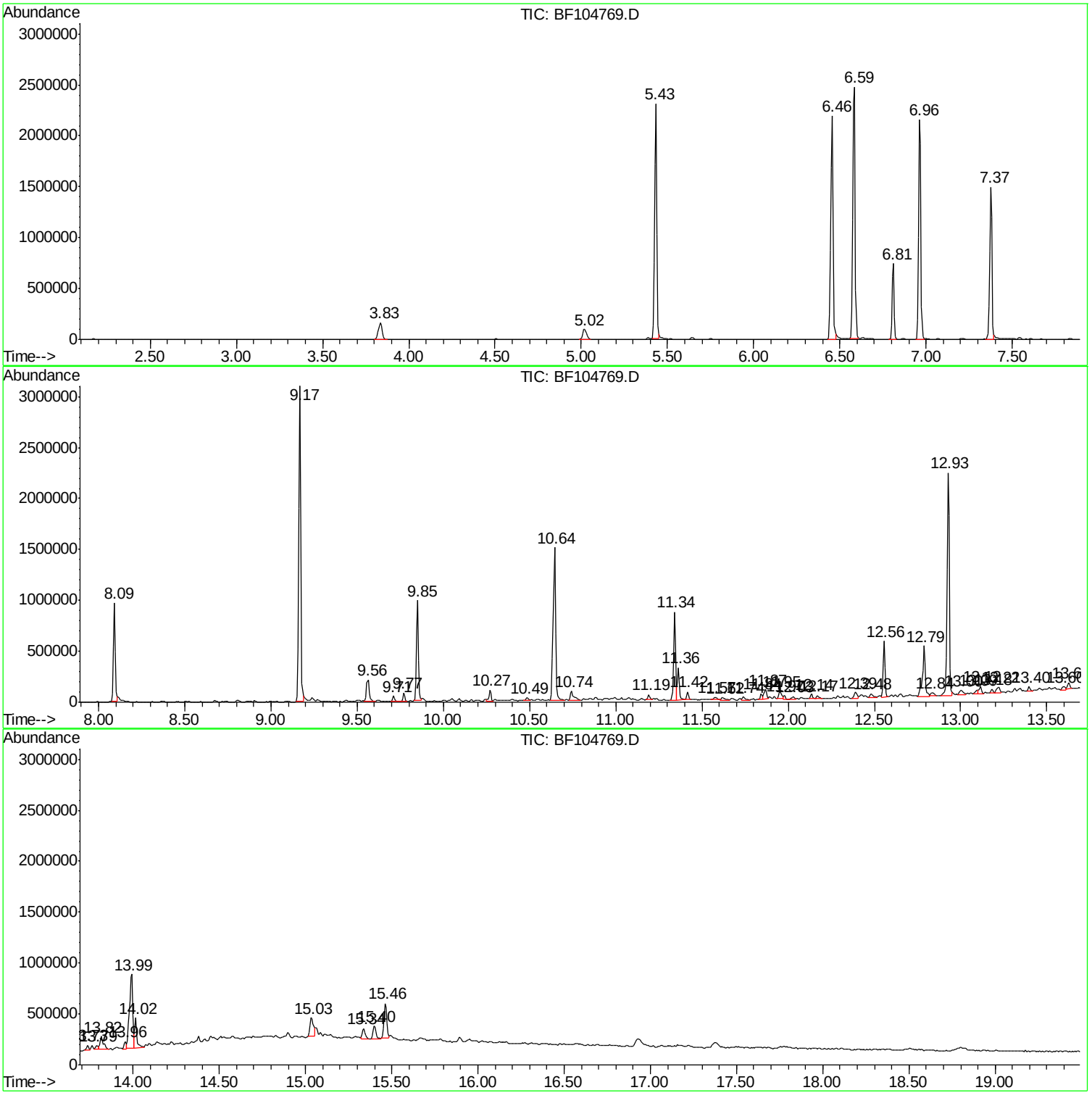
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

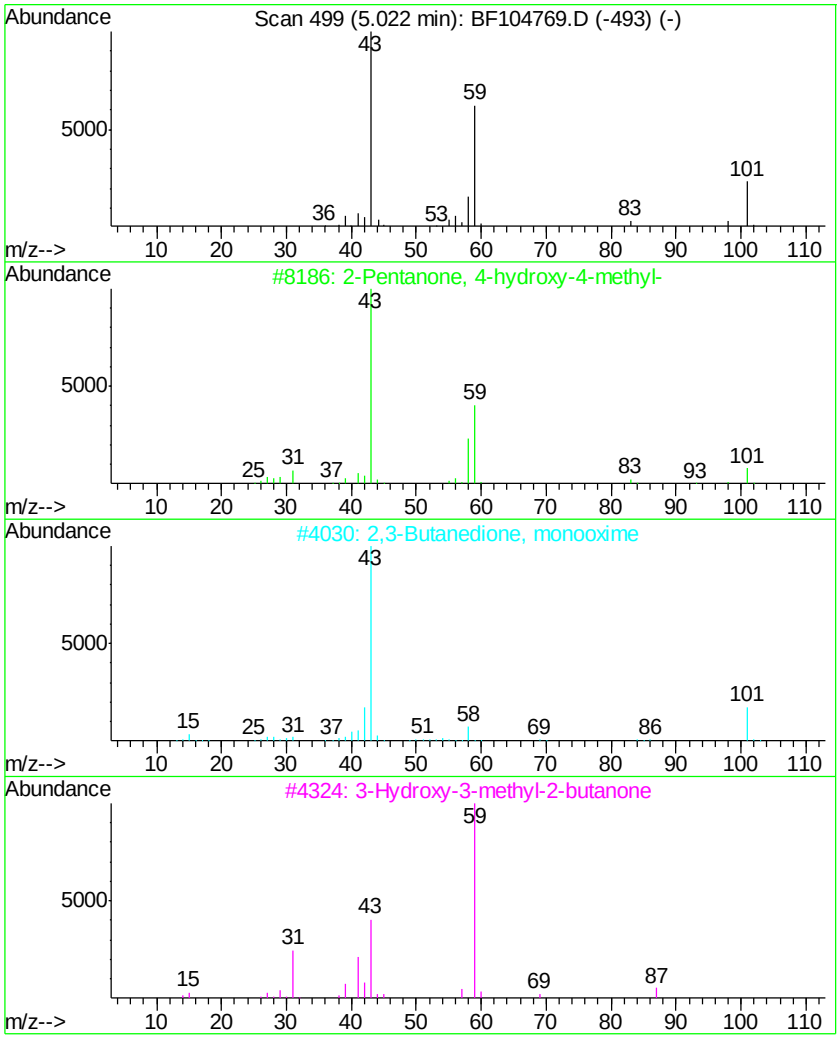
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

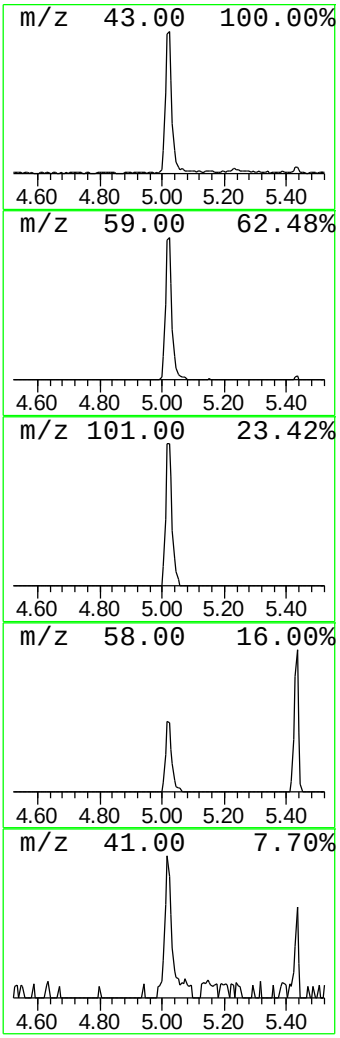
**********

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.02 | 4.59 ng | 141669 | 1,4-Dichlorobenzene-d4 | 6.81 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 50 |
| 2 | 2,3-Butanedione, monooxime | 101 | C4H7NO2 | 000057-71-6 | 35 |
| 3 | 3-Hydroxy-3-methyl-2-butanone | 102 | C5H10O2 | 000115-22-0 | 12 |
| 4 | 5-Hexen-2-one | 98 | C6H10O | 000109-49-9 | 9 |
| 5 | (+-)-4-Amino-4,5-dihydro-2(3H)-f... | 101 | C4H7NO2 | 016504-58-8 | 9 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

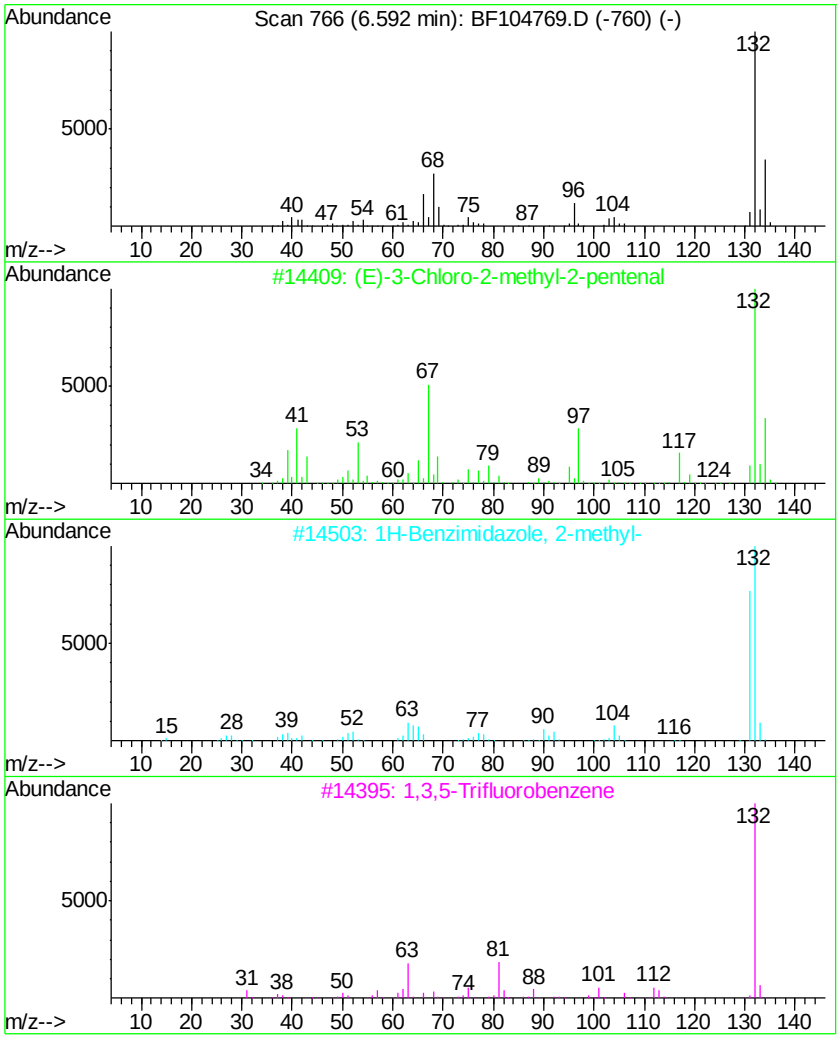
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

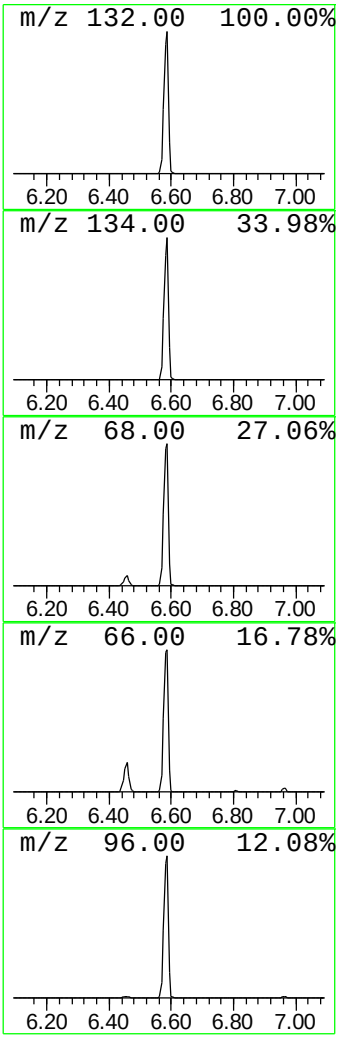
**********

Peak Number 3 unknown6.59 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.59 | 77.42 ng | 2389100 | 1,4-Dichlorobenzene-d4 | 6.81 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO | 031357-76-3 | 17 |
| 2 | 1H-Benzimidazole, 2-methyl- | 132 | C8H8N2 | 000615-15-6 | 10 |
| 3 | 1,3,5-Trifluorobenzene | 132 | C6H3F3 | 000372-38-3 | 9 |
| 4 | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2 | 022291-85-6 | 9 |
| 5 | 5-Fluoro-2-chloropyrimidine | 132 | C4H2ClFN2 | 062802-42-0 | 9 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

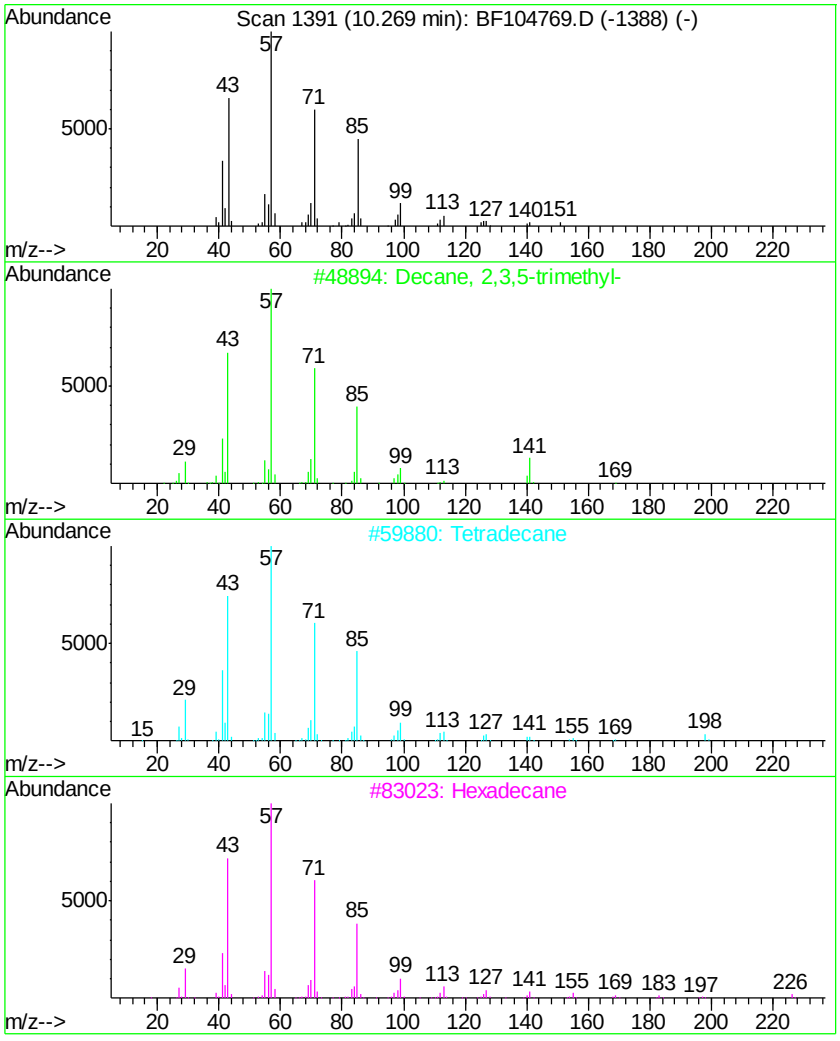
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

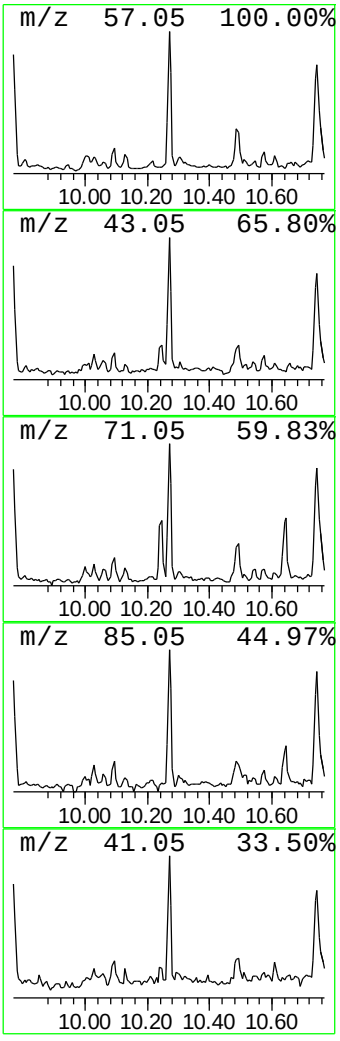
**********

Peak Number 5 Decane, 2,3,5-trimethyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.27 | 2.10 ng | 87756 | Acenaphthene-d10 | 9.85 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Decane, 2,3,5-trimethyl- | 184 | C13H28 | 062238-11-3 | 90 |
| 2 | Tetradecane | 198 | C14H30 | 000629-59-4 | 87 |
| 3 | Hexadecane | 226 | C16H34 | 000544-76-3 | 78 |
| 4 | Heptadecane | 240 | C17H36 | 000629-78-7 | 78 |
| 5 | Tridecane | 184 | C13H28 | 000629-50-5 | 59 |



# Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

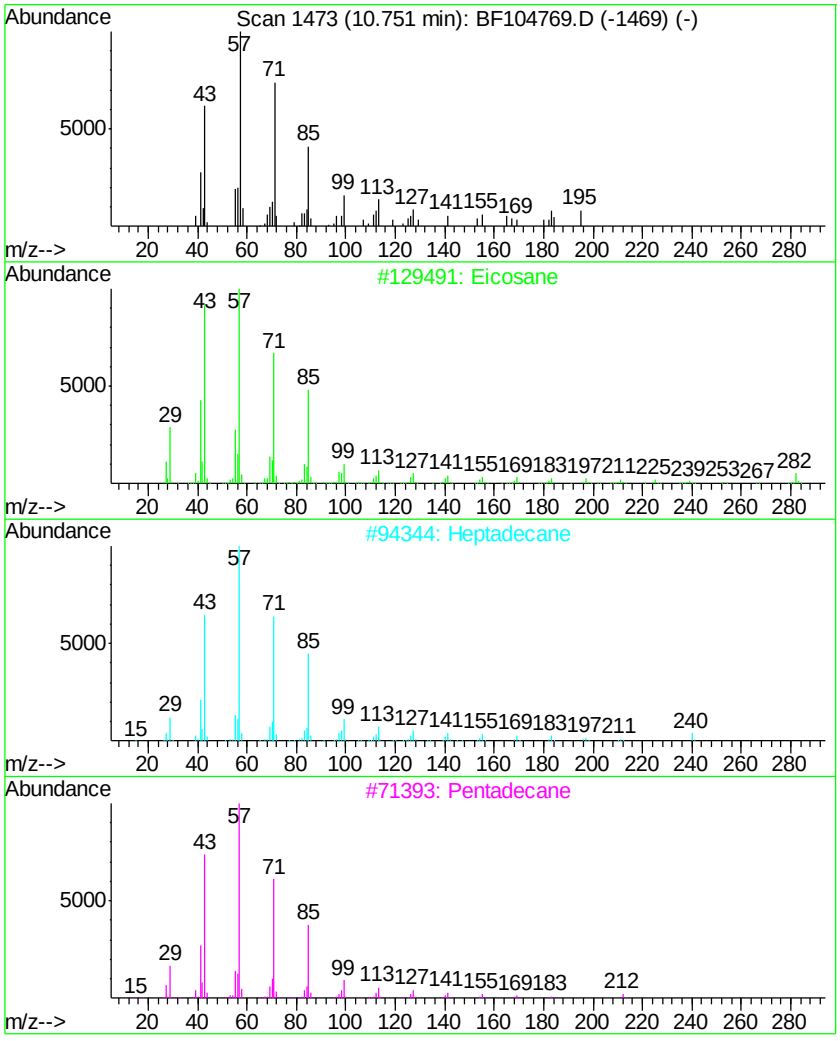
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

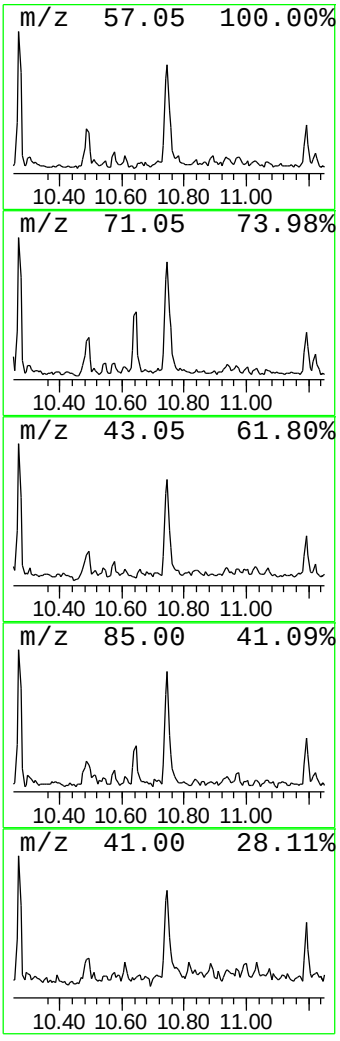
**********

Peak Number 6 Eicosane Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.75 | 3.35 ng | 123543 | Phenanthrene-d10 | 11.34 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Eicosane | 282 | C20H42 | 000112-95-8 | 86 |
| 2 | Heptadecane | 240 | C17H36 | 000629-78-7 | 86 |
| 3 | Pentadecane | 212 | C15H32 | 000629-62-9 | 86 |
| 4 | Hexadecane | 226 | C16H34 | 000544-76-3 | 86 |
| 5 | Tritetracontane | 605 | C43H88 | 007098-21-7 | 86 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

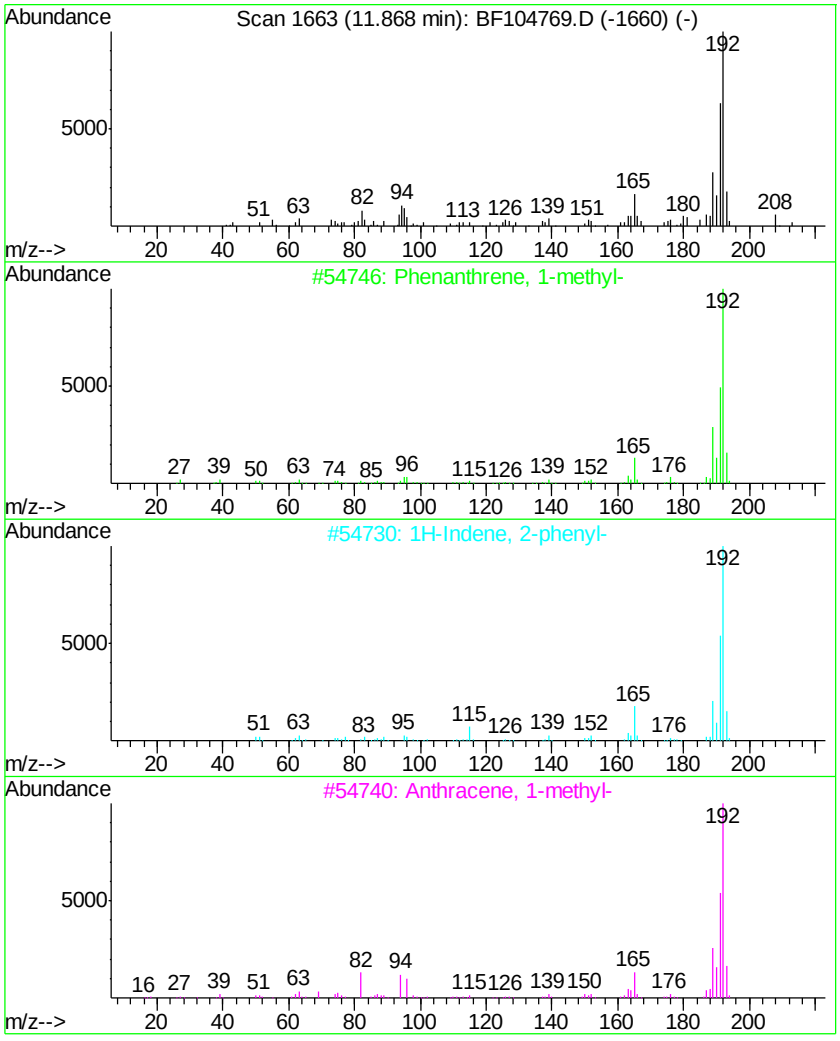
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

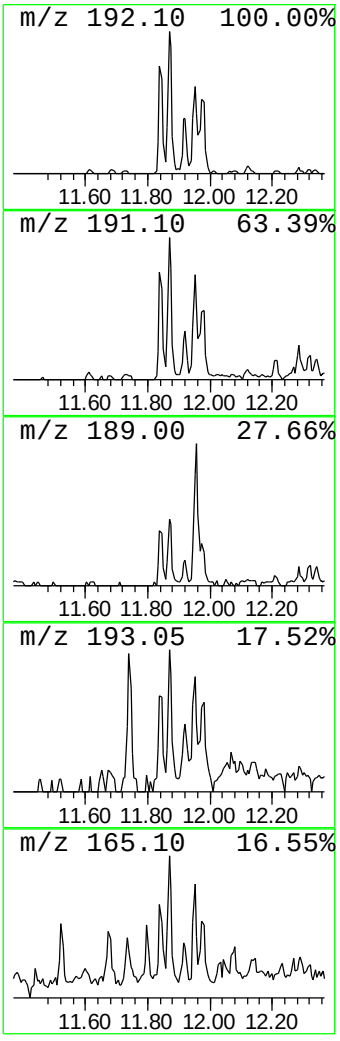
**********

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

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.87 | 2.77 ng | 102100 | Phenanthrene-d10 | 11.34 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 95 |
| 2 | 1H-Indene, 2-phenyl- | 192 | C15H12 | 004505-48-0 | 94 |
| 3 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 93 |
| 4 | Anthracene, 9-methyl- | 192 | C15H12 | 000779-02-2 | 91 |
| 5 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 90 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

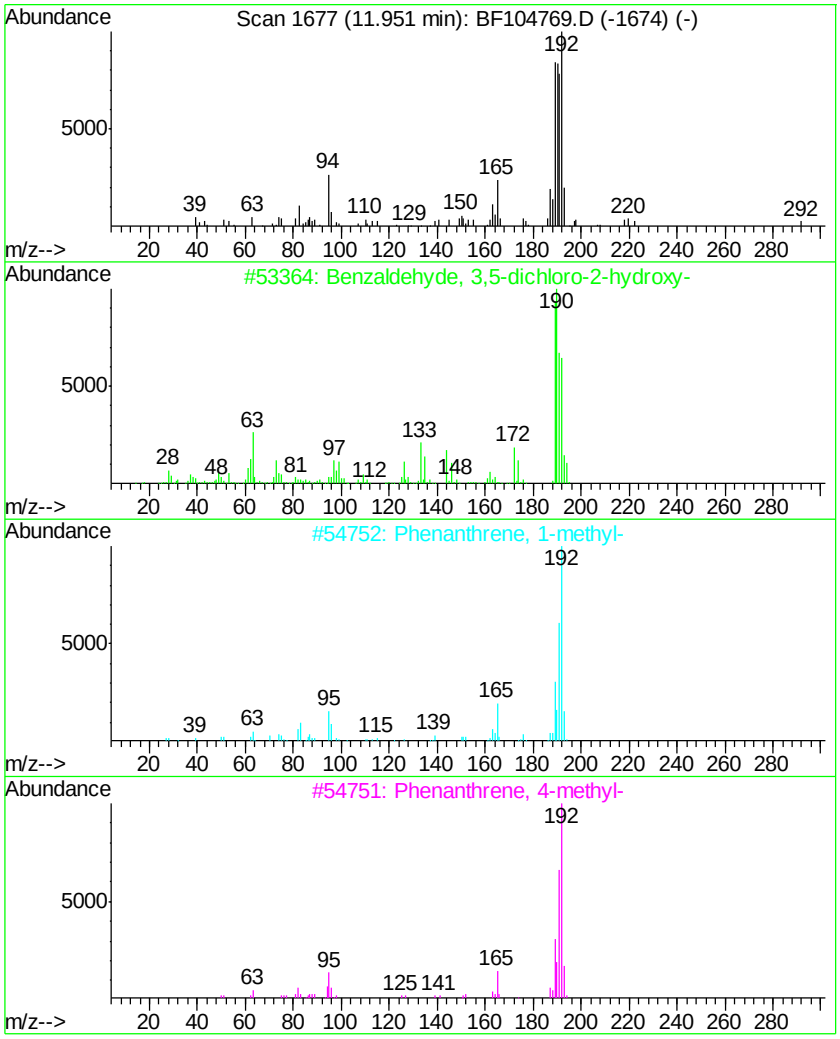
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

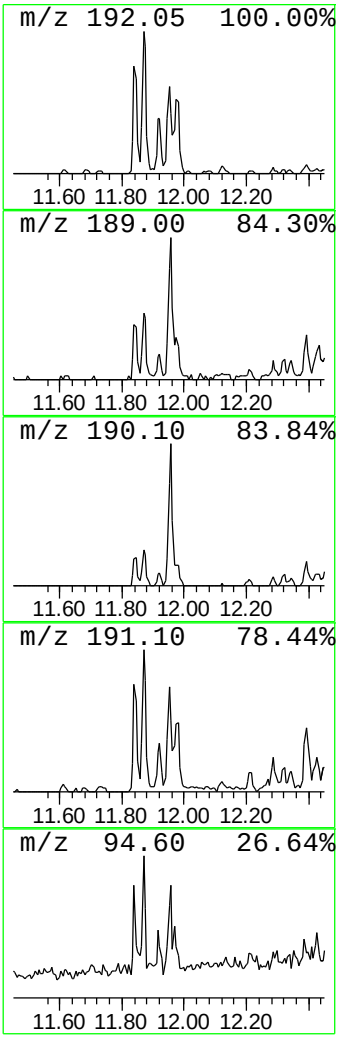
**********

Peak Number 8 Benzaldehyde, 3,5-dichloro-... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.95 | 2.18 ng | 80515 | Phenanthrene-d10 | 11.34 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzaldehyde, 3,5-dichloro-2-hyd... | 190 | C7H4Cl2O2 | 000090-60-8 | 58 |
| 2 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 50 |
| 3 | Phenanthrene, 4-methyl- | 192 | C15H12 | 000832-64-4 | 50 |
| 4 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 50 |
| 5 | 6H-Cyclobuta[jk]phenanthrene | 190 | C15H10 | 083469-43-6 | 43 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

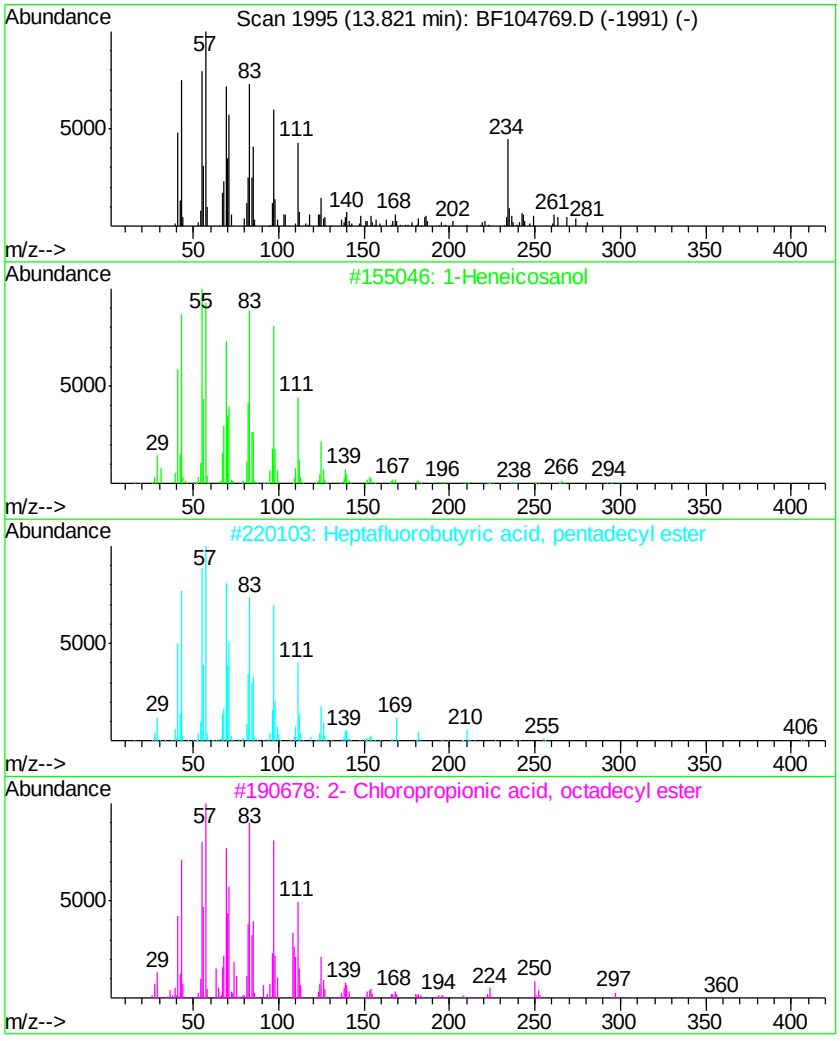
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

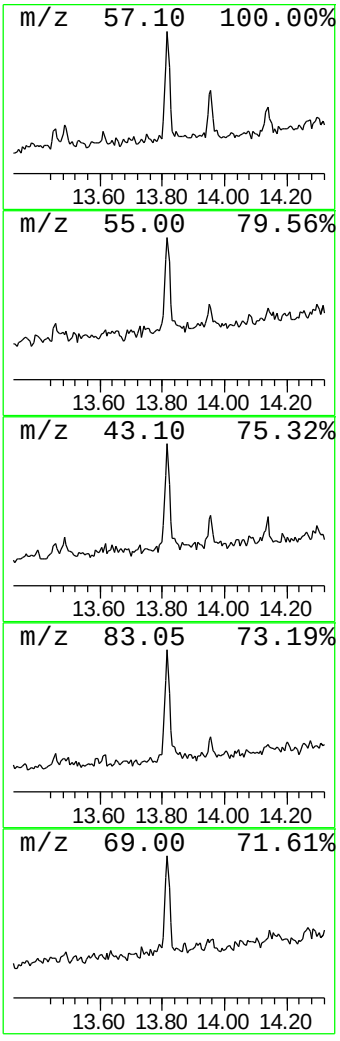
**********

Peak Number 9 1-Heneicosanol Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.82 | 3.82 ng | 179532 | Chrysene-d12 | 13.99 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Heneicosanol | 312 | C21H44O | 015594-90-8 | 93 |
| 2 | Heptafluorobutyric acid, pentade... | 424 | C19H31F7O2 | 959261-23-5 | 93 |
| 3 | 2- Chloropropionic acid, octadec... | 360 | C21H41ClO2 | 088104-31-8 | 90 |
| 4 | 1-Heptacosanol | 396 | C27H56O | 002004-39-9 | 87 |
| 5 | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 83 |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

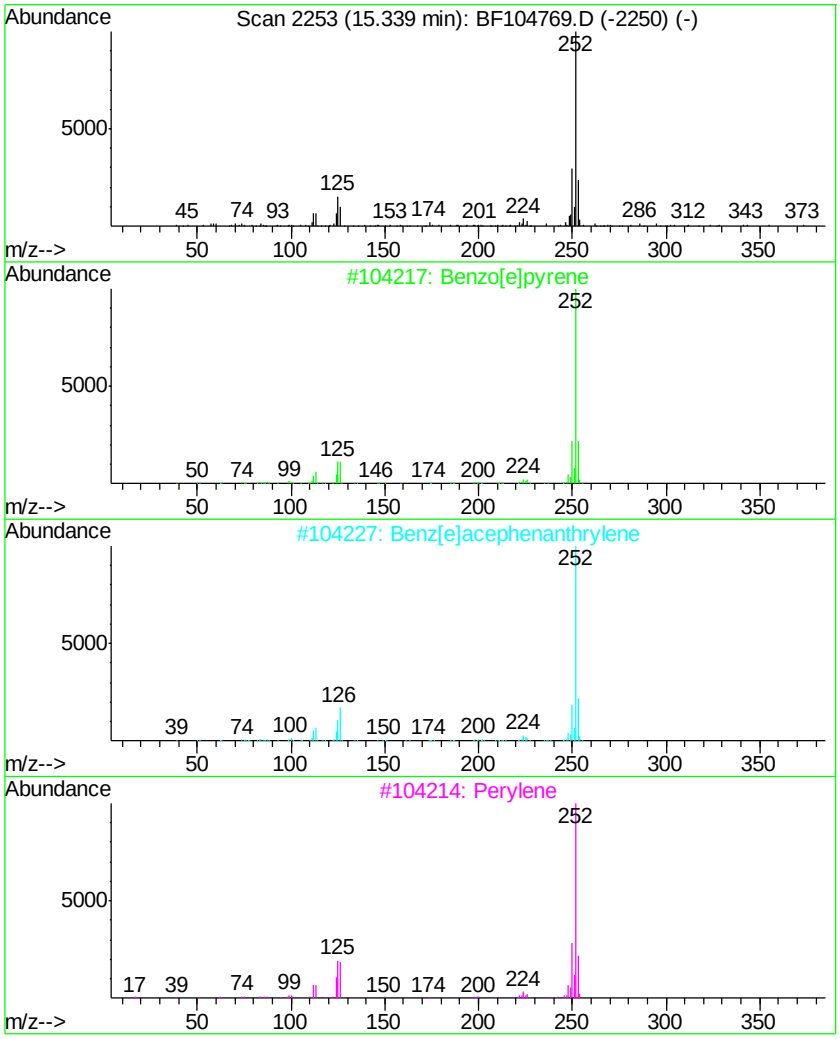
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

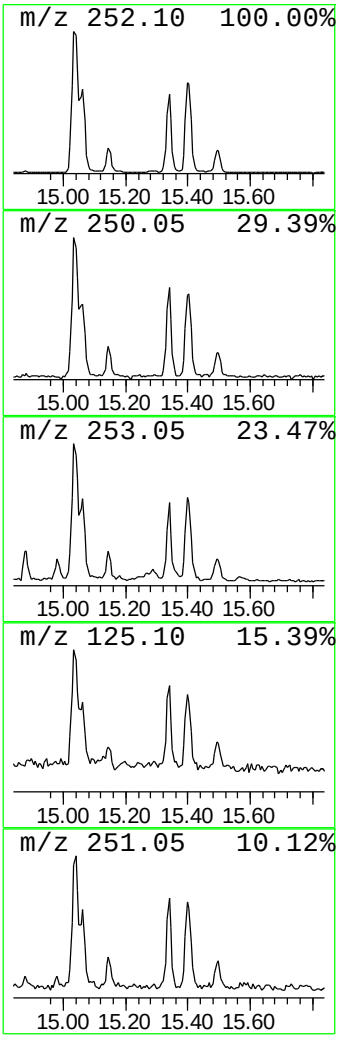
**********

Peak Number 10 Benzo[e]pyrene Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.34 | 5.93 ng | 125625 | Perylene-d12 | 15.46 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 93 |
| 2 | Benz[e]acephenanthrylene | 252 | C20H12 | 000205-99-2 | 90 |
| 3 | Perylene | 252 | C20H12 | 000198-55-0 | 90 |
| 4 | Benzo[k]fluoranthene | 252 | C20H12 | 000207-08-9 | 89 |
| 5 | Benzo[a]pyrene | 252 | C20H12 | 000050-32-8 | 81 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042418\
Data File : BF104769.D
Acq On : 24 Apr 2018 21:01
Operator : JU/SJ
Sample : J2508-13
Misc :
ALS Vial : 20 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** TP-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.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... | 5.02 | 4.6 | ng | 141669 | 1 | 6.81 | 617182 | 20.0 |
| unknown6.59 | 6.59 | 77.4 | ng | 2389100 | 1 | 6.81 | 617182 | 20.0 |
| Decane, 2,3,5-tri... | 10.27 | 2.1 | ng | 87756 | 3 | 9.85 | 836053 | 20.0 |
| Eicosane | 10.75 | 3.4 | ng | 123543 | 4 | 11.34 | 737990 | 20.0 |
| Phenanthrene, 1-m... | 11.87 | 2.8 | ng | 102100 | 4 | 11.34 | 737990 | 20.0 |
| Benzaldehyde, 3,5... | 11.95 | 2.2 | ng | 80515 | 4 | 11.34 | 737990 | 20.0 |
| 1-Heneicosanol | 13.82 | 3.8 | ng | 179532 | 5 | 13.99 | 941064 | 20.0 |
| Benzo[e]pyrene | 15.34 | 5.9 | ng | 125625 | 6 | 15.46 | 423919 | 20.0 |