LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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 | 5.440 | 566 | 570 | 582 | rVB | 1227479 | 1141275 | 81.41% | 7.619% |
| 2 | 6.457 | 738 | 743 | 758 | rVV | 1321208 | 1260380 | 89.91% | 8.414% |
| 3 | 6.593 | 762 | 766 | 774 | rBV | 1449615 | 1291202 | 92.11% | 8.619% |
| 4 | 6.651 | 774 | 776 | 782 | rVB2 | 18021 | 17863 | 1.27% | 0.119% |
| 5 | 6.822 | 802 | 805 | 813 | rBV | 393463 | 370804 | 26.45% | 2.475% |
| 6 | 6.981 | 828 | 832 | 843 | rVB | 1158009 | 975953 | 69.62% | 6.515% |
| 7 | 7.381 | 896 | 900 | 919 | rVB | 792458 | 762966 | 54.43% | 5.093% |
| 8 | 8.104 | 1019 | 1023 | 1033 | rBV | 478232 | 492439 | 35.13% | 3.287% |
| 9 | 9.175 | 1201 | 1205 | 1216 | rBV | 1676710 | 1401862 | 100.00% | 9.358% |
| 10 | 9.257 | 1216 | 1219 | 1230 | rVB5 | 25690 | 37101 | 2.65% | 0.248% |
| 11 | 9.563 | 1268 | 1271 | 1280 | rBV | 114965 | 115362 | 8.23% | 0.770% |
| 12 | 9.716 | 1294 | 1297 | 1301 | rVB | 20019 | 21037 | 1.50% | 0.140% |
| 13 | 9.851 | 1315 | 1320 | 1324 | rBV | 583501 | 523804 | 37.36% | 3.497% |
| 14 | 9.881 | 1324 | 1325 | 1331 | rVB3 | 21722 | 22251 | 1.59% | 0.149% |
| 15 | 10.057 | 1351 | 1355 | 1360 | rBV2 | 19391 | 26972 | 1.92% | 0.180% |
| 16 | 10.398 | 1410 | 1413 | 1418 | rVB | 27396 | 31082 | 2.22% | 0.207% |
| 17 | 10.639 | 1450 | 1454 | 1465 | rBV | 745579 | 775727 | 55.34% | 5.178% |
| 18 | 10.769 | 1472 | 1476 | 1482 | rVB3 | 35165 | 55210 | 3.94% | 0.369% |
| 19 | 11.233 | 1553 | 1555 | 1562 | rVB2 | 34051 | 34145 | 2.44% | 0.228% |
| 20 | 11.333 | 1568 | 1572 | 1574 | rBV | 484643 | 429412 | 30.63% | 2.867% |
| 21 | 11.357 | 1574 | 1576 | 1582 | rVV | 203045 | 185609 | 13.24% | 1.239% |
| 22 | 11.410 | 1582 | 1585 | 1587 | rVV | 59693 | 49157 | 3.51% | 0.328% |
| 23 | 11.628 | 1620 | 1622 | 1626 | rVB5 | 24868 | 21466 | 1.53% | 0.143% |
| 24 | 11.839 | 1655 | 1658 | 1659 | rBV | 27441 | 26252 | 1.87% | 0.175% |
| 25 | 11.863 | 1659 | 1662 | 1668 | rVV3 | 162445 | 168626 | 12.03% | 1.126% |
| 26 | 11.945 | 1674 | 1676 | 1679 | rVB | 51687 | 45191 | 3.22% | 0.302% |
| 27 | 12.128 | 1705 | 1707 | 1710 | rBV | 25389 | 23371 | 1.67% | 0.156% |
| 28 | 12.545 | 1775 | 1778 | 1782 | rVB | 441204 | 366219 | 26.12% | 2.445% |
| 29 | 12.586 | 1783 | 1785 | 1791 | rVB4 | 53032 | 51736 | 3.69% | 0.345% |
| 30 | 12.645 | 1792 | 1795 | 1798 | rBV3 | 62737 | 70595 | 5.04% | 0.471% |
| 31 | 12.775 | 1812 | 1817 | 1820 | rBV | 401513 | 395196 | 28.19% | 2.638% |
| 32 | 12.916 | 1837 | 1841 | 1844 | rBV | 1042367 | 905614 | 64.60% | 6.045% |
| 33 | 13.104 | 1870 | 1873 | 1876 | rVB | 71138 | 75625 | 5.39% | 0.505% |
| 34 | 13.169 | 1882 | 1884 | 1886 | rVB | 49089 | 36279 | 2.59% | 0.242% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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 | 13.969 | 2015 | 2020 | 2023 | rBV2 | 454931 | 603484 | 43.05% | 4.029% |
| 36 | 13.992 | 2023 | 2024 | 2030 | rVB | 207001 | 165875 | 11.83% | 1.107% |
| 37 | 14.486 | 2105 | 2108 | 2109 | rBV2 | 64949 | 73125 | 5.22% | 0.488% |
| 38 | 14.998 | 2191 | 2195 | 2203 | rBV | 269610 | 483431 | 34.48% | 3.227% |
| 39 | 15.298 | 2243 | 2246 | 2251 | rVB | 151147 | 194784 | 13.89% | 1.300% |
| 40 | 15.357 | 2252 | 2256 | 2262 | rVV | 194100 | 292406 | 20.86% | 1.952% |
| 41 | 15.416 | 2262 | 2266 | 2277 | rVB2 | 306377 | 539189 | 38.46% | 3.599% |
| 42 | 16.527 | 2452 | 2455 | 2462 | rVB3 | 83183 | 126852 | 9.05% | 0.847% |
| 43 | 16.845 | 2506 | 2509 | 2516 | rVB2 | 95915 | 173265 | 12.36% | 1.157% |
| 44 | 17.280 | 2579 | 2583 | 2592 | rVB | 59827 | 119918 | 8.55% | 0.801% |

Sum of corrected areas: 14980112

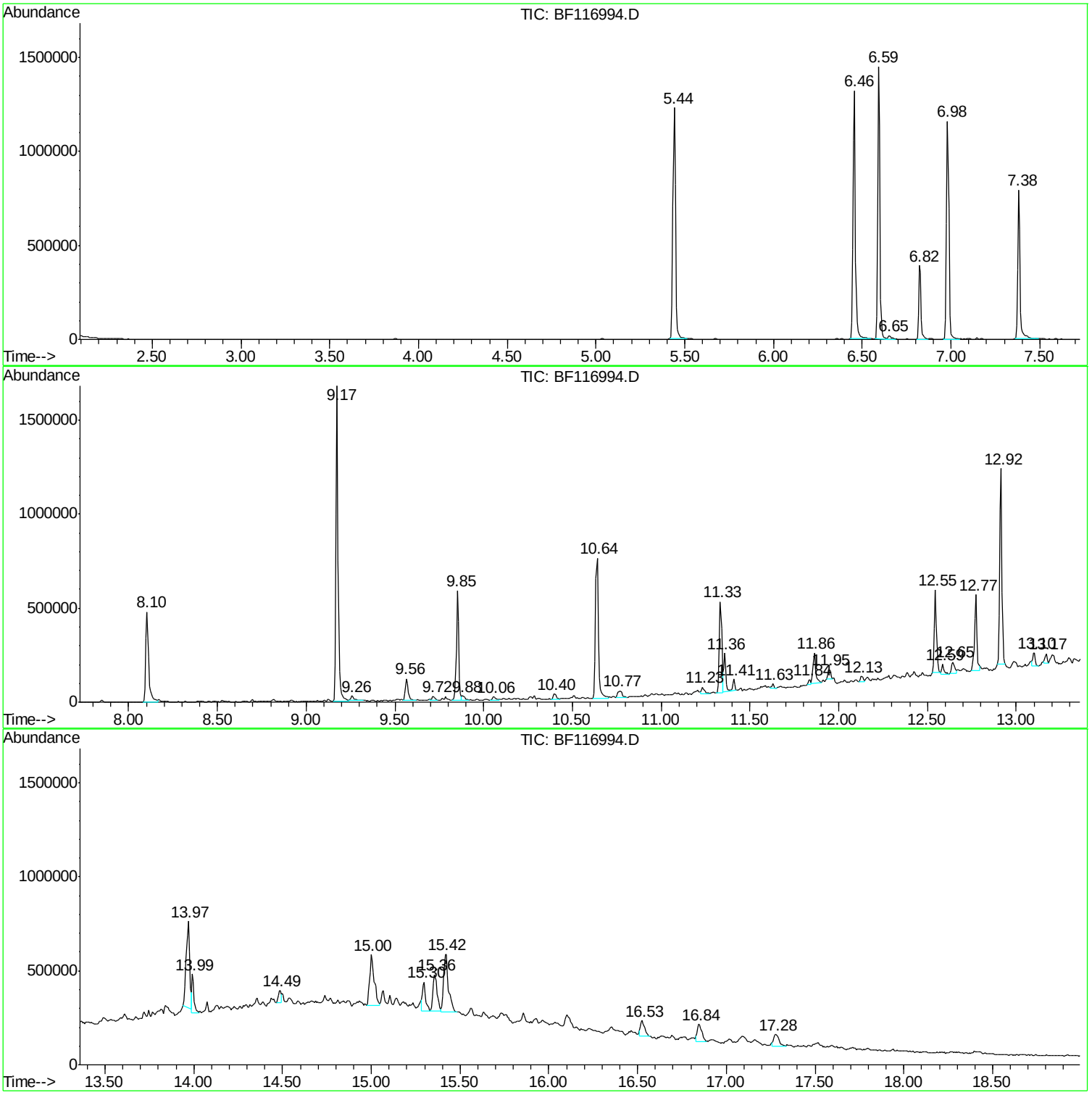
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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

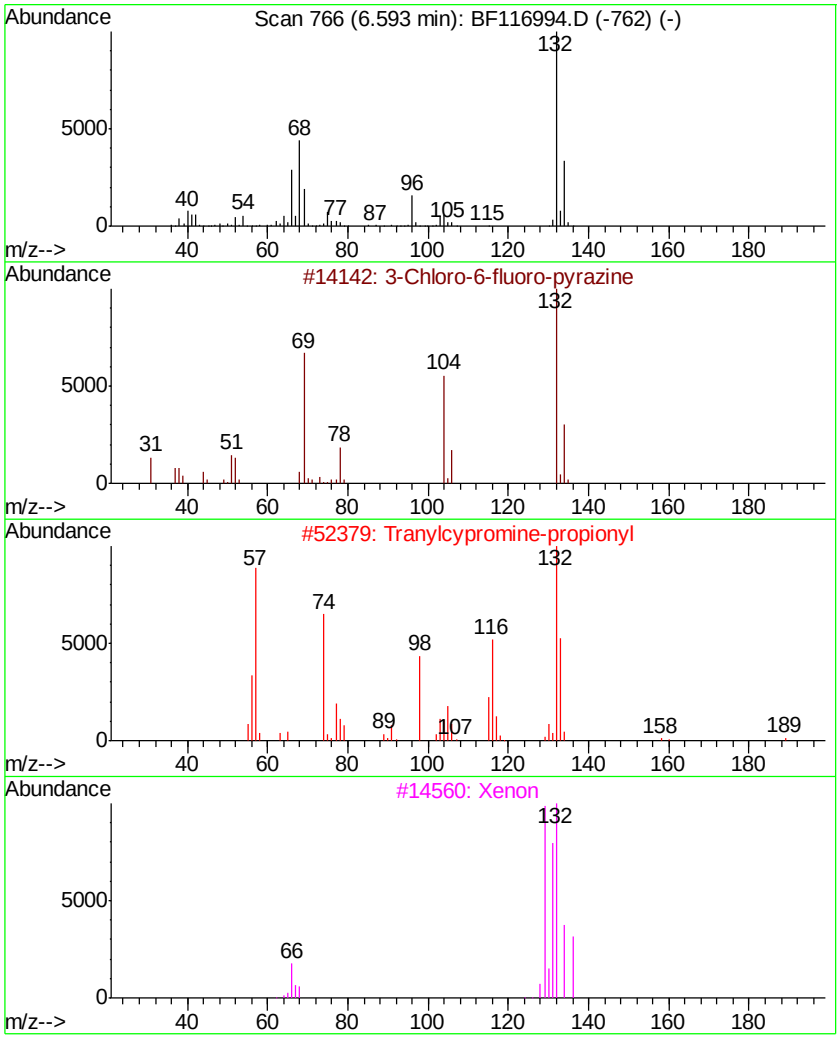
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

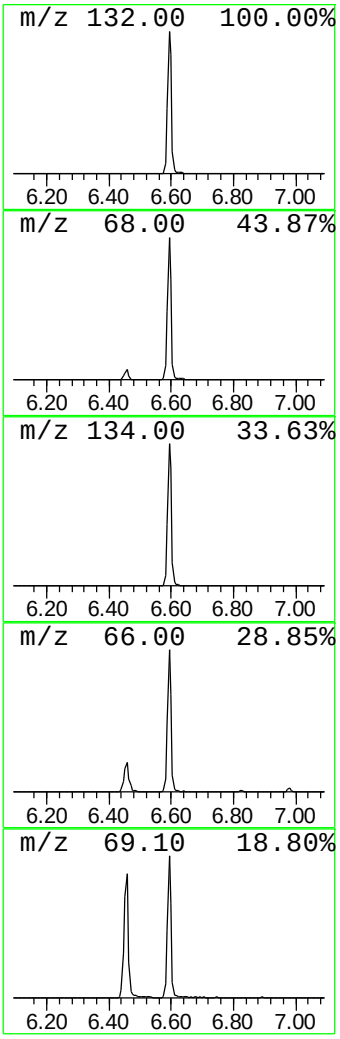
**********

Peak Number 1 unknown6.59 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.59 | 69.64 ng | 1291200 | 1,4-Dichlorobenzene-d4 | 6.82 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Chloro-6-fluoro-pyrazine | 132 | C4H2ClFN2 | 1000146-10-7 | 16 |
| 2 | Tranylcypromine-propionyl | 189 | C12H15NO | 1000123-86-3 | 10 |
| 3 | Xenon | 132 | Xe | 007440-63-3 | 9 |
| 4 | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12 | 028749-81-7 | 9 |
| 5 | Benzene, 1-ethenyl-3,5-dimethyl- | 132 | C10H12 | 005379-20-4 | 9 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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

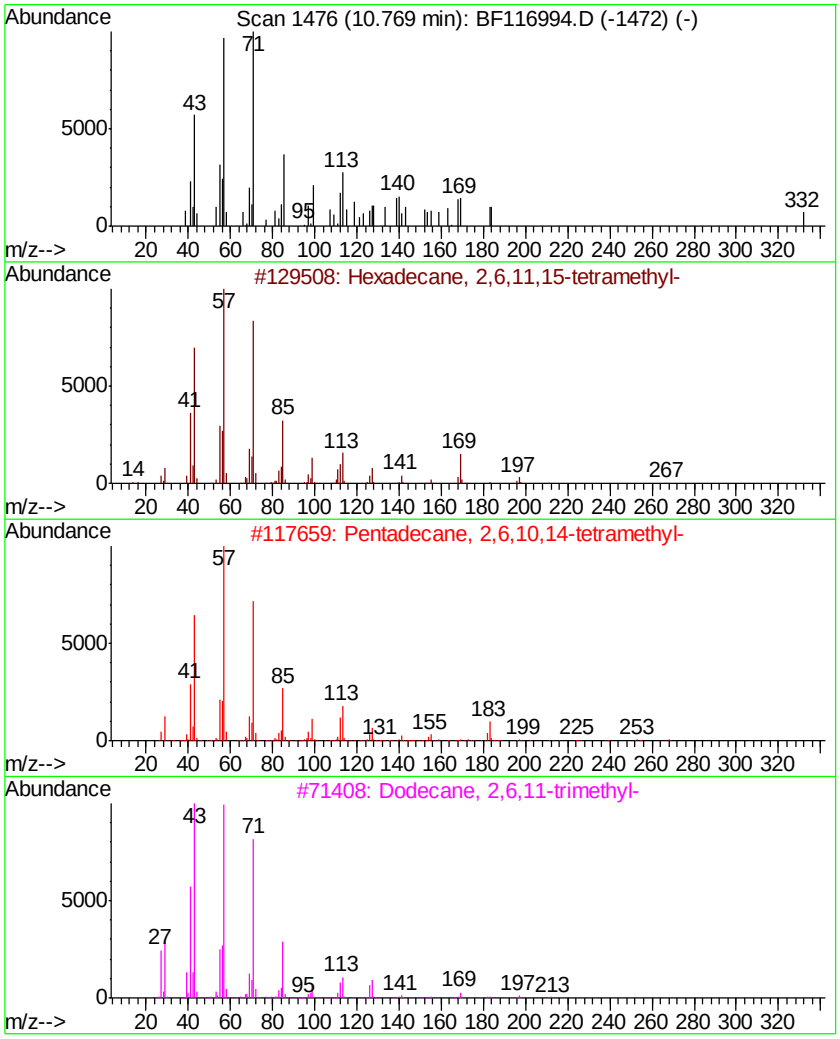
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

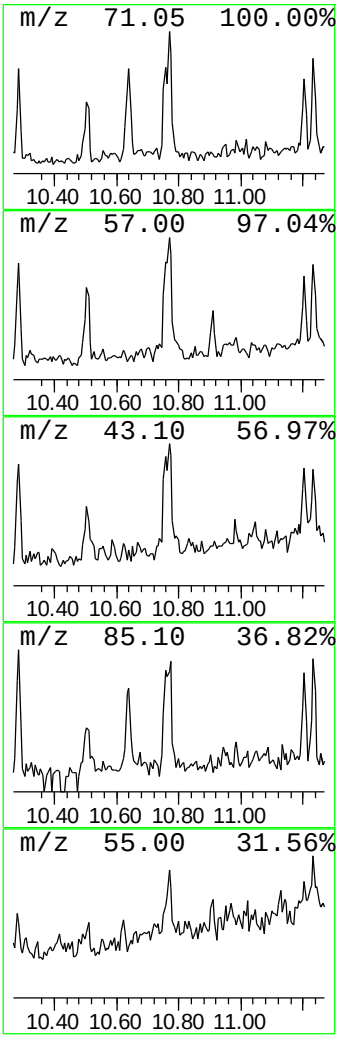
**********

Peak Number 3 Hexadecane, 2,6,11,15-tetra... Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.77 | 2.57 ng | 55210 | Phenanthrene-d10 | 11.33 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexadecane, 2,6,11,15-tetramethyl- | 282 | C20H42 | 000504-44-9 | 64 |
| 2 | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40 | 001921-70-6 | 50 |
| 3 | Dodecane, 2,6,11-trimethyl- | 212 | C15H32 | 031295-56-4 | 47 |
| 4 | Sulfurous acid, 2-ethylhexyl hep... | 432 | C25H52O3S | 1000309-20-0 | 43 |
| 5 | Octane, 3,6-dimethyl- | 142 | C10H22 | 015869-94-0 | 43 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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

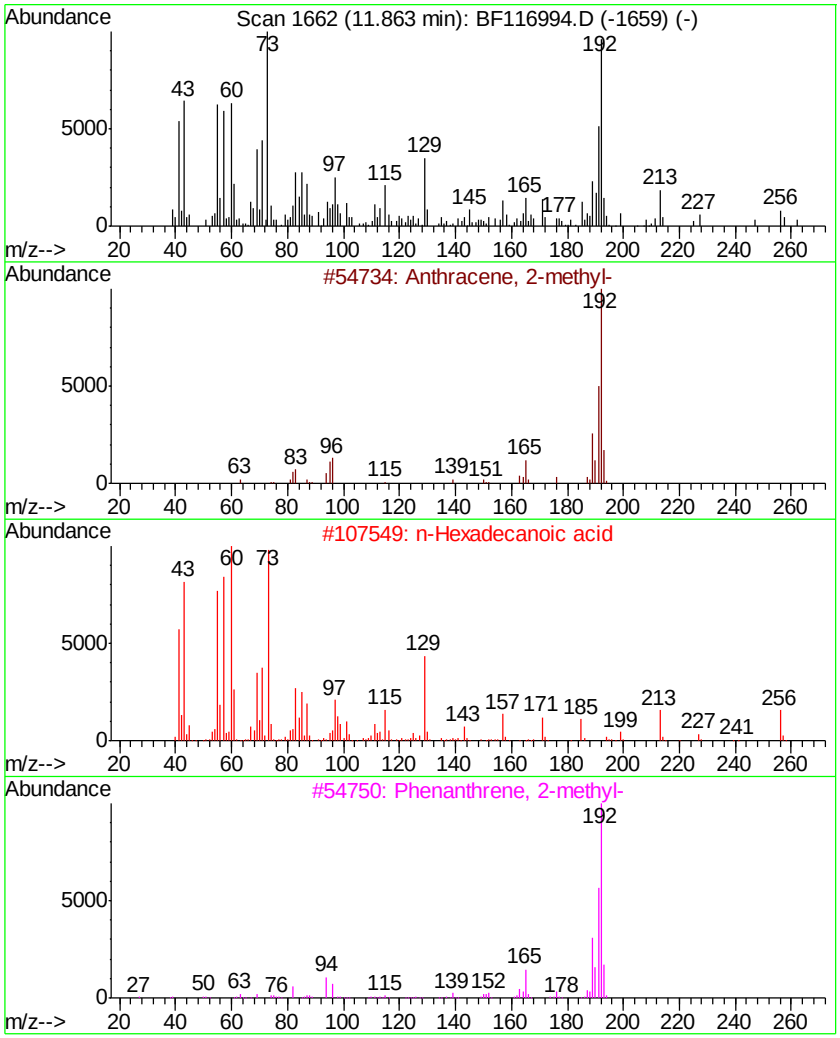
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

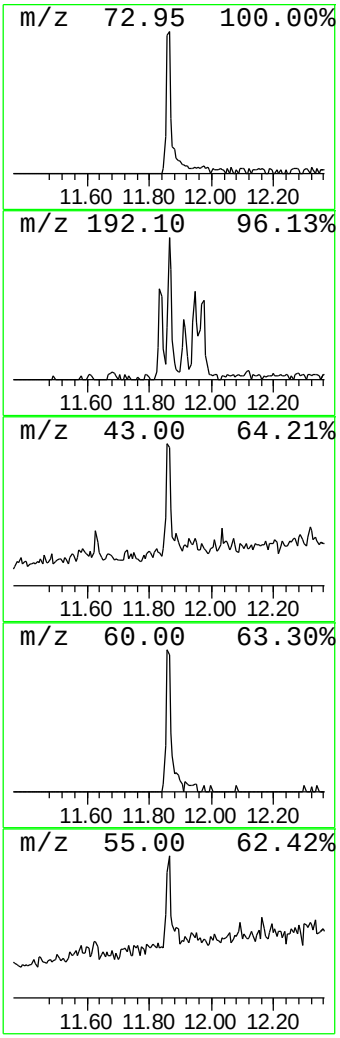
**********

Peak Number 4 Anthracene, 2-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.86 | 7.85 ng | 168626 | Phenanthrene-d10 | 11.33 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 80 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 72 |
| 3 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 51 |
| 4 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 47 |
| 5 | Anthracene, 9-methyl- | 192 | C15H12 | 000779-02-2 | 47 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

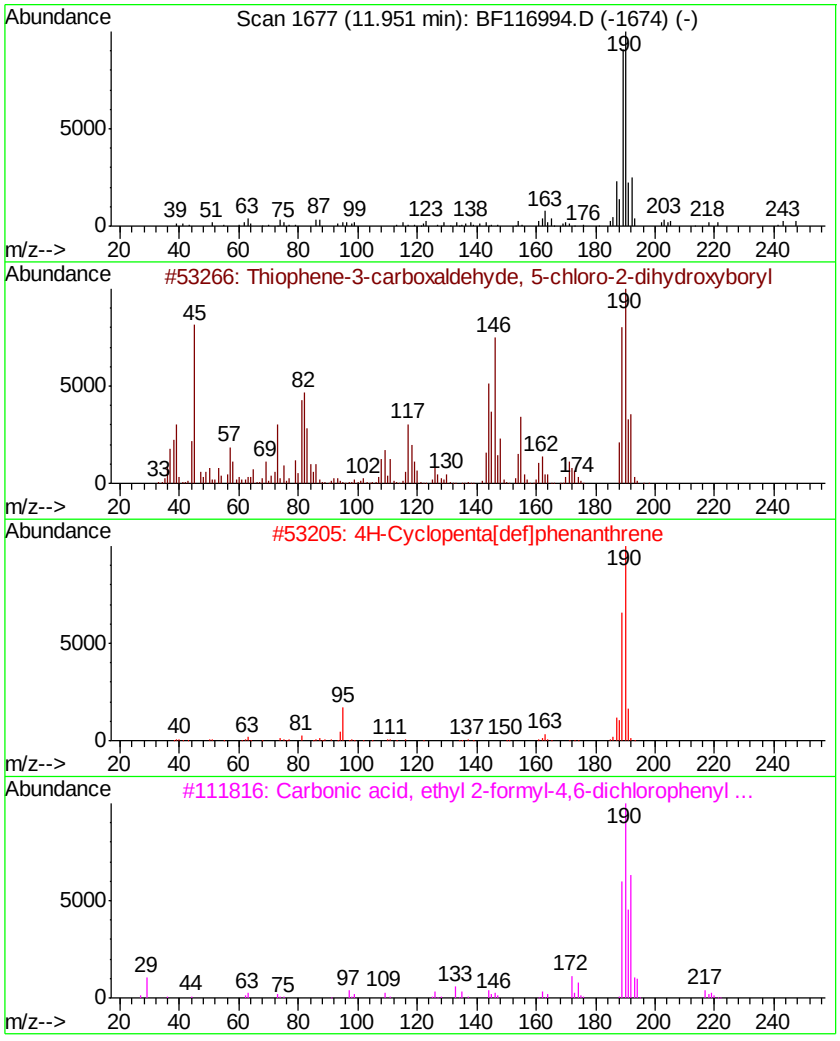
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

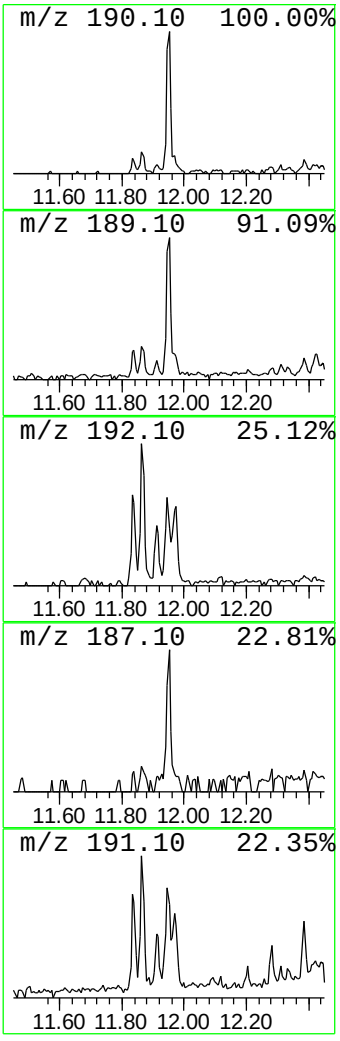
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 Thiophene-3-carboxaldehyde,... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.95 | 2.10 ng | 45191 | Phenanthrene-d10 | 11.33 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Thiophene-3-carboxaldehyde, 5-ch... | 190 | C5H4BClO3S | 036155-87-0 | 72 |
| 2 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 68 |
| 3 | Carbonic acid, ethyl 2-formyl-4,... | 262 | C10H8Cl2O4 | 1000331-34-4 | 50 |
| 4 | 6H-Cyclobuta[jk]phenanthrene | 190 | C15H10 | 083469-43-6 | 42 |
| 5 | Benzaldehyde, 3,5-dichloro-2-hyd... | 190 | C7H4Cl2O2 | 000090-60-8 | 42 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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

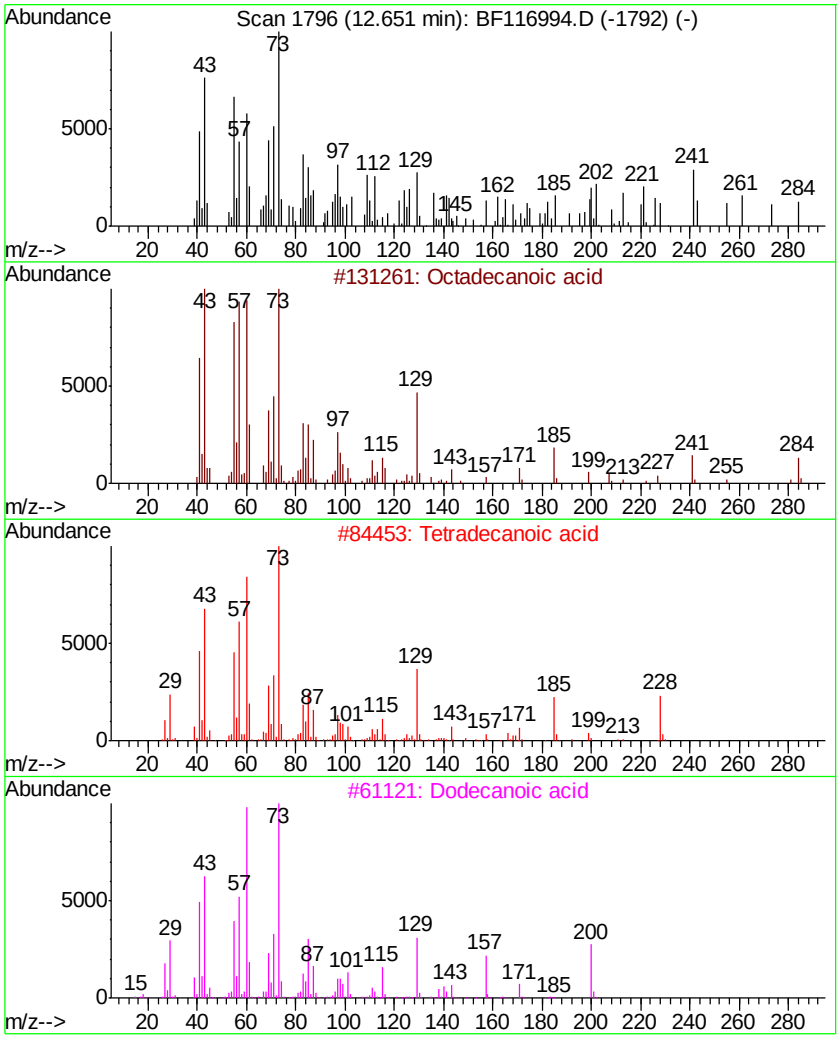
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

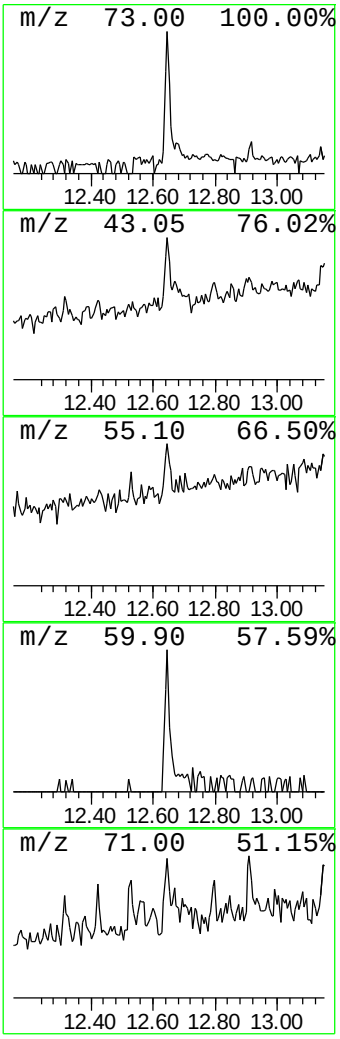
**********

Peak Number 6 Octadecanoic acid Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.65 | 3.29 ng | 70595 | Phenanthrene-d10 | 11.33 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 64 |
| 2 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 52 |
| 3 | Dodecanoic acid | 200 | C12H24O2 | 000143-07-7 | 49 |
| 4 | Lactose | 342 | C12H22O11 | 000063-42-3 | 38 |
| 5 | n-Decanoic acid | 172 | C10H20O2 | 000334-48-5 | 38 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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

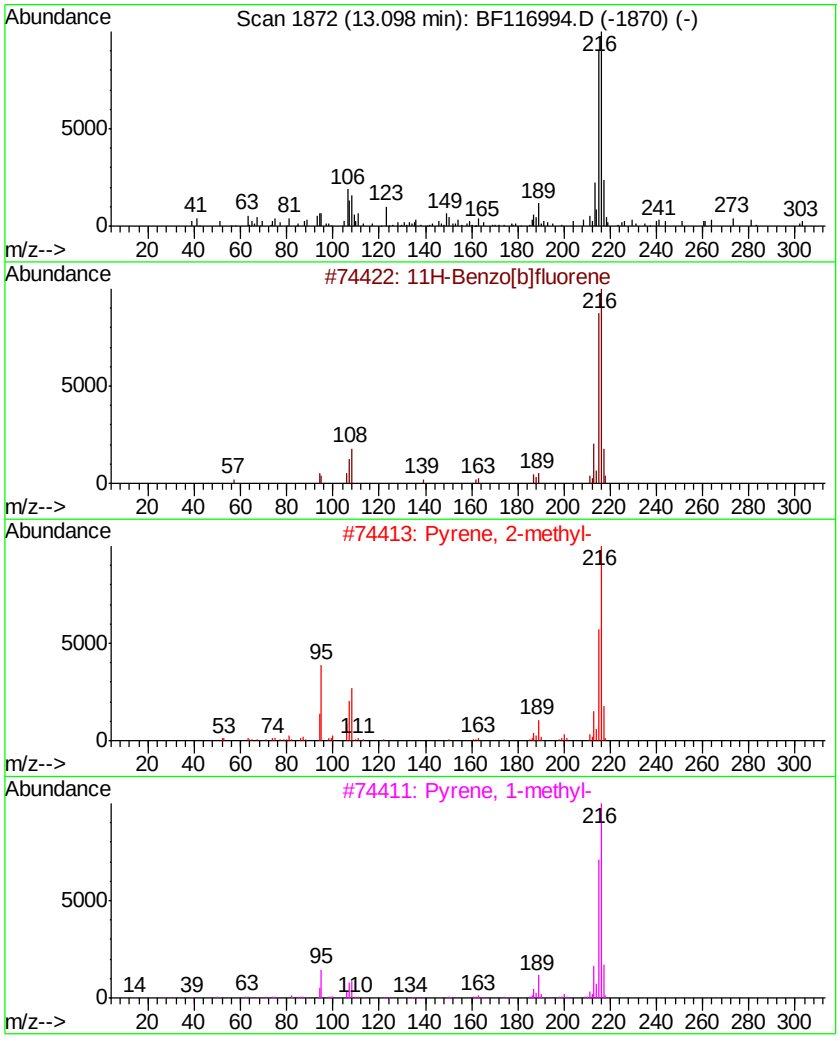
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

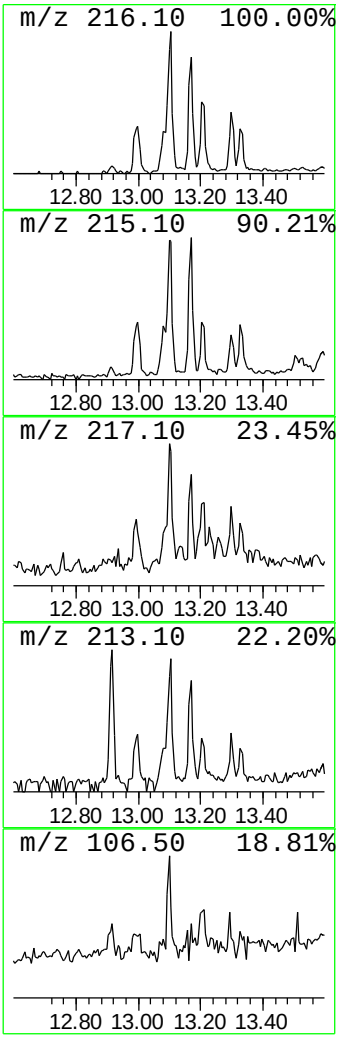
**********

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.10 | 2.51 ng | 75625 | Chrysene-d12 | 13.97 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 91 |
| 2 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 86 |
| 3 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 86 |
| 4 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 83 |
| 5 | 7H-Benzo[c]fluorene | 216 | C17H12 | 000205-12-9 | 74 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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

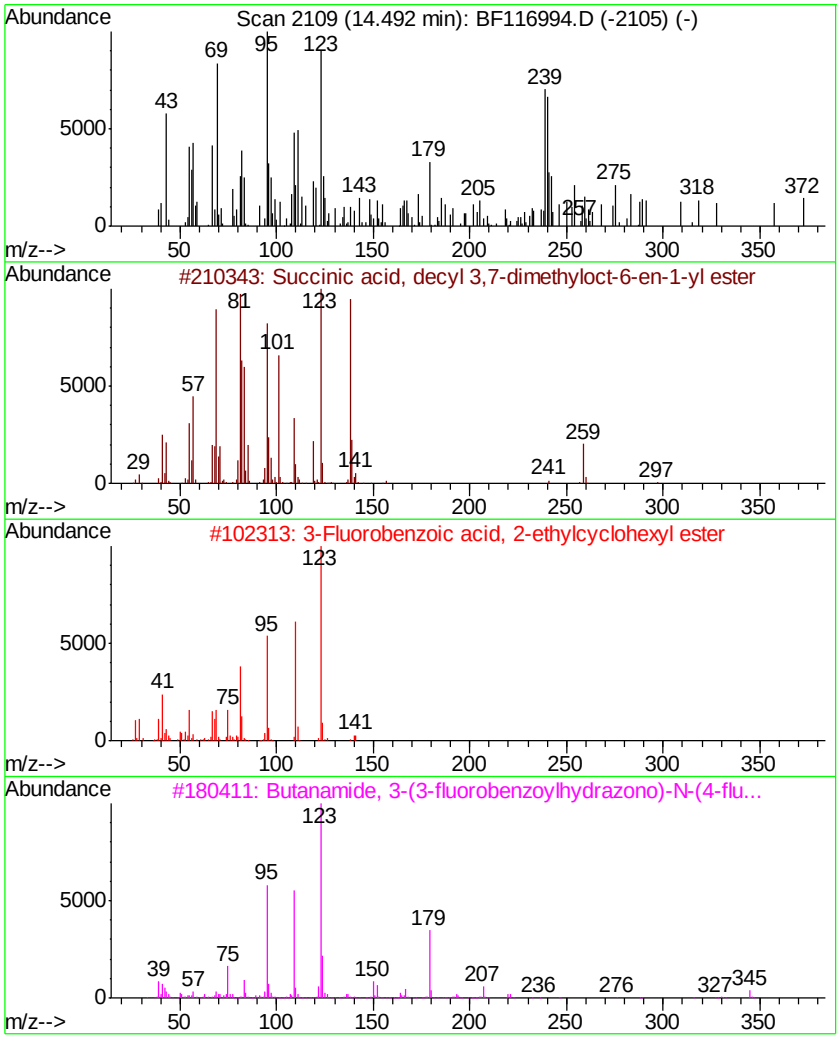
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

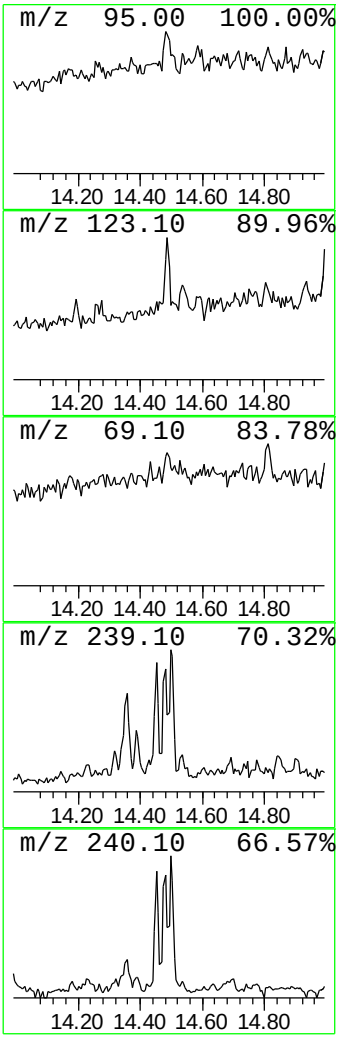
**********

Peak Number 8 unknown14.49 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.49 | 2.42 ng | 73125 | Chrysene-d12 | 13.97 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Succinic acid, decyl 3,7-dimethyl... | 396 | C24H44O4 | 1000353-34-4 | 47 |
| 2 | 3-Fluorobenzoic acid, 2-ethylcyc... | 250 | C15H19FO2 | 1000279-04-9 | 43 |
| 3 | Butanamide, 3-(3-fluorobenzoylhy... | 345 | C18H17F2N3O2 | 1000260-11-6 | 35 |
| 4 | trans-Cyclohexanecarboxylic acid... | 240 | C15H28O2 | 038289-32-6 | 27 |
| 5 | Benzamide, 4-fluoro-N-(5-chloro-... | 250 | C12H8ClFN2O | 327991-18-4 | 25 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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

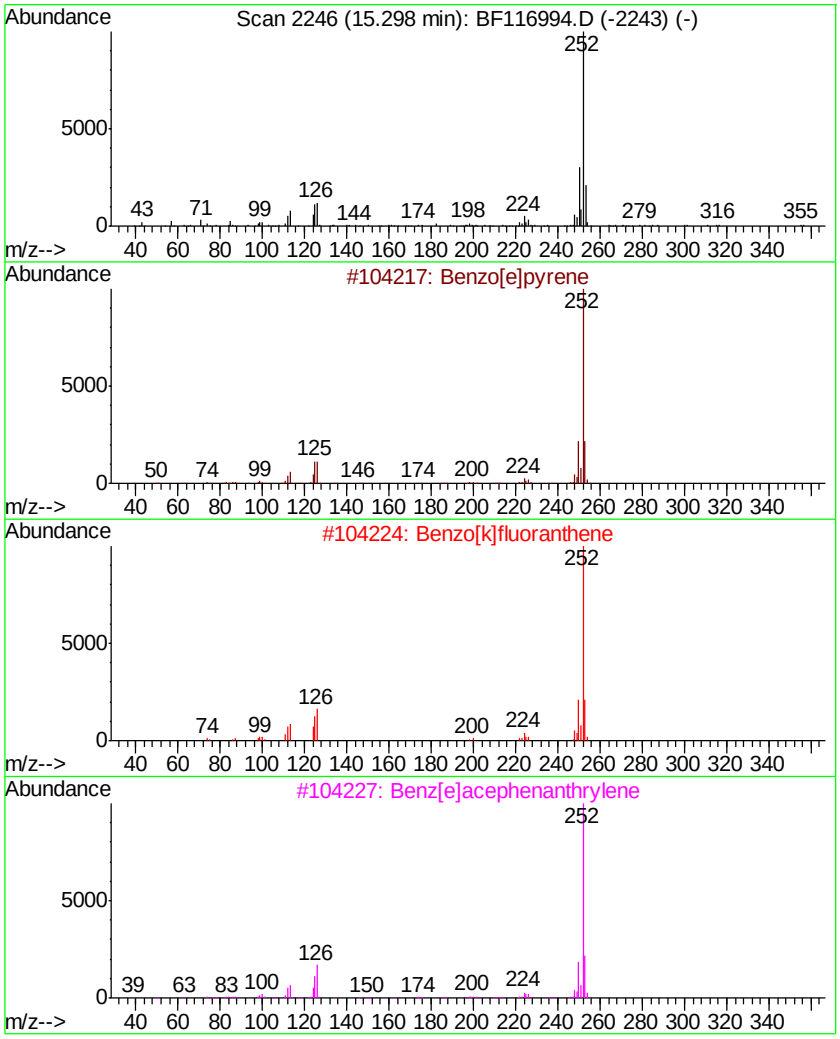
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

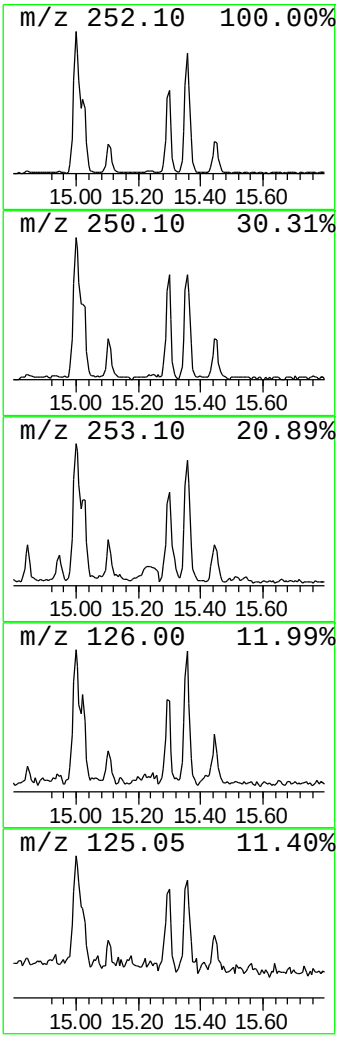
**********

Peak Number 9 Benzo[e]pyrene Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.30 | 7.23 ng | 194784 | Perylene-d12 | 15.42 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 98 |
| 2 | Benzo[k]fluoranthene | 252 | C20H12 | 000207-08-9 | 98 |
| 3 | Benz[e]acephenanthrylene | 252 | C20H12 | 000205-99-2 | 96 |
| 4 | Benzo[a]pyrene | 252 | C20H12 | 000050-32-8 | 96 |
| 5 | Perylene | 252 | C20H12 | 000198-55-0 | 93 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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

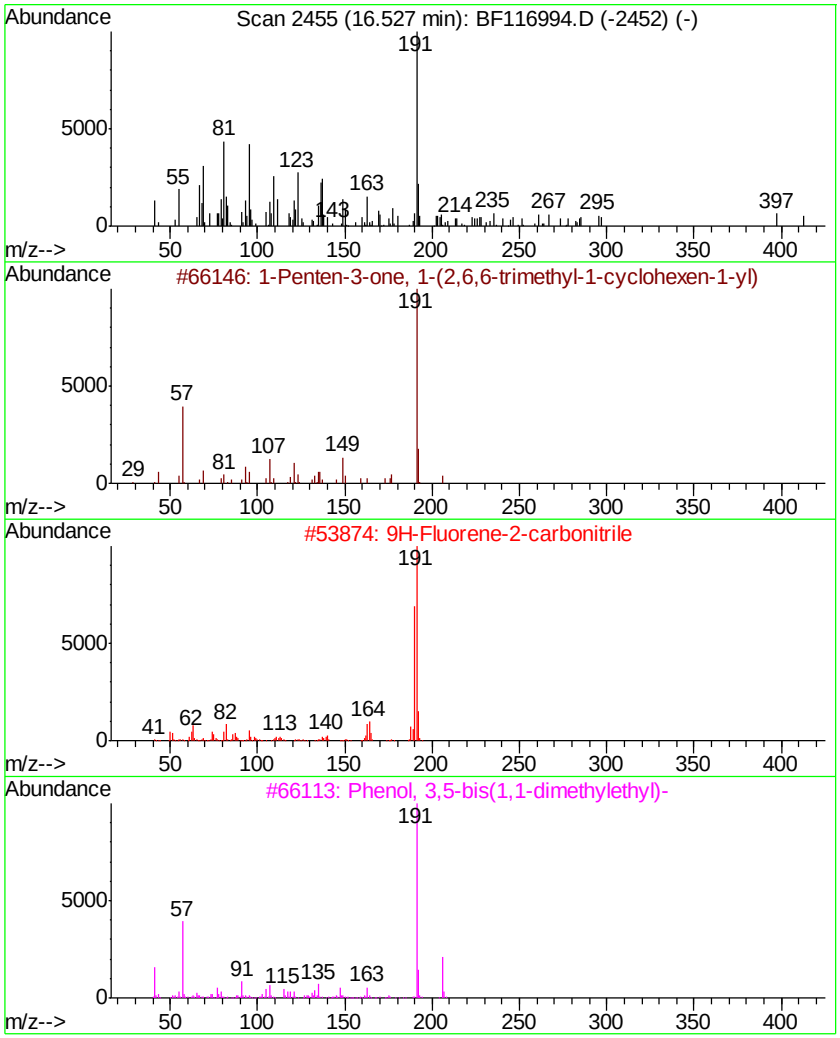
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

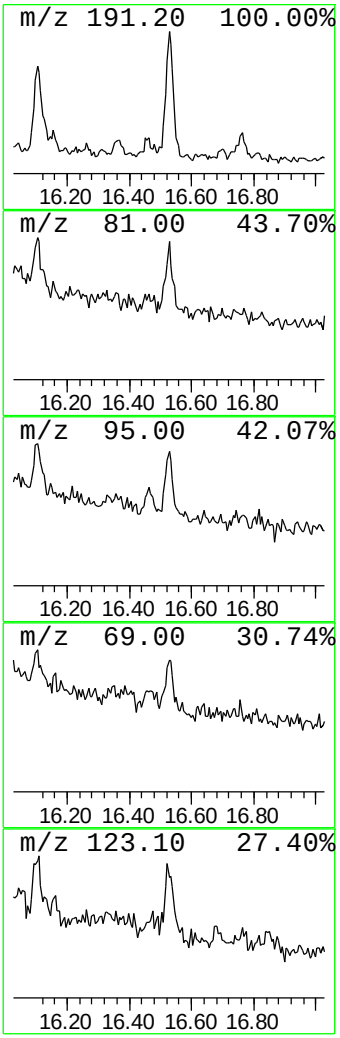
**********

Peak Number 10 unknown16.53 Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.53 | 4.71 ng | 126852 | Perylene-d12 | 15.42 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Penten-3-one, 1-(2,6,6-trimeth... | 206 | C14H22O | 000127-43-5 | 35 |
| 2 | 9H-Fluorene-2-carbonitrile | 191 | C14H9N | 002523-48-0 | 35 |
| 3 | Phenol, 3,5-bis(1,1-dimethylethyl)- | 206 | C14H22O | 001138-52-9 | 27 |
| 4 | 2-Thiazolemethanol, .alpha.-phenyl- | 191 | C10H9NOS | 000879-52-7 | 27 |
| 5 | Amiphenazole | 191 | C9H9N3S | 000490-55-1 | 25 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF091219\
Data File : BF116994.D
Acq On : 12 Sep 2019 16:15
Operator : HP/JU
Sample : K4852-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** BUS-3(0-3.5)

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 |
| unknown6.59 | 6.59 | 69.6 | ng | 1291200 | 1 | 6.82 | 370804 | 20.0 |
| Hexadecane, 2,6,1... | 10.77 | 2.6 | ng | 55210 | 4 | 11.33 | 429412 | 20.0 |
| Anthracene, 2-met... | 11.86 | 7.8 | ng | 168626 | 4 | 11.33 | 429412 | 20.0 |
| Thiophene-3-carbo... | 11.95 | 2.1 | ng | 45191 | 4 | 11.33 | 429412 | 20.0 |
| Octadecanoic acid | 12.65 | 3.3 | ng | 70595 | 4 | 11.33 | 429412 | 20.0 |
| 11H-Benzo[b]fluorene | 13.10 | 2.5 | ng | 75625 | 5 | 13.97 | 603484 | 20.0 |
| unknown14.49 | 14.49 | 2.4 | ng | 73125 | 5 | 13.97 | 603484 | 20.0 |
| Benzo[e]pyrene | 15.30 | 7.2 | ng | 194784 | 6 | 15.42 | 539189 | 20.0 |
| unknown16.53 | 16.53 | 4.7 | ng | 126852 | 6 | 15.42 | 539189 | 20.0 |