LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

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-BF040622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127827.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 2.328 | 3 | 7 | 13 | rVB | 263545 | 329646 | 7.13% | 0.911% |
| 2 | 5.204 | 493 | 496 | 503 | rVB | 35610 | 47974 | 1.04% | 0.133% |
| 3 | 5.581 | 552 | 560 | 563 | rBV | 3696947 | 3773032 | 81.56% | 10.427% |
| 4 | 6.575 | 722 | 729 | 732 | rBV | 3426425 | 3879738 | 83.86% | 10.722% |
| 5 | 6.945 | 789 | 792 | 796 | rVB | 1084721 | 950904 | 20.55% | 2.628% |
| 6 | 7.516 | 882 | 889 | 892 | rBV | 2303763 | 2669728 | 57.71% | 7.378% |
| 7 | 8.234 | 1006 | 1011 | 1014 | rBV | 1321458 | 1242200 | 26.85% | 3.433% |
| 8 | 9.310 | 1188 | 1194 | 1197 | rBV | 4959234 | 4626320 | 100.00% | 12.785% |
| 9 | 9.992 | 1301 | 1310 | 1313 | rBV | 1726370 | 1467734 | 31.73% | 4.056% |
| 10 | 10.780 | 1438 | 1444 | 1448 | rBV | 3022603 | 3120422 | 67.45% | 8.624% |
| 11 | 11.480 | 1559 | 1563 | 1566 | rBV | 1636872 | 1516248 | 32.77% | 4.190% |
| 12 | 11.504 | 1566 | 1567 | 1571 | rVV | 194353 | 121959 | 2.64% | 0.337% |
| 13 | 11.557 | 1571 | 1576 | 1578 | rVV | 92579 | 81622 | 1.76% | 0.226% |
| 14 | 11.710 | 1598 | 1602 | 1609 | rBV3 | 30417 | 55064 | 1.19% | 0.152% |
| 15 | 11.998 | 1642 | 1651 | 1657 | rBV2 | 207243 | 315856 | 6.83% | 0.873% |
| 16 | 12.098 | 1664 | 1668 | 1670 | rBV2 | 146491 | 135688 | 2.93% | 0.375% |
| 17 | 12.533 | 1739 | 1742 | 1745 | rBV2 | 50618 | 60537 | 1.31% | 0.167% |
| 18 | 12.574 | 1745 | 1749 | 1754 | rVB4 | 38215 | 60386 | 1.31% | 0.167% |
| 19 | 12.621 | 1754 | 1757 | 1762 | rBV | 117273 | 116688 | 2.52% | 0.322% |
| 20 | 12.698 | 1766 | 1770 | 1774 | rBV | 1135341 | 972970 | 21.03% | 2.689% |
| 21 | 12.927 | 1803 | 1809 | 1815 | rBV2 | 877854 | 1018291 | 22.01% | 2.814% |
| 22 | 12.974 | 1815 | 1817 | 1822 | rVB2 | 44882 | 52682 | 1.14% | 0.146% |
| 23 | 13.074 | 1826 | 1834 | 1837 | rBV | 4470544 | 4272348 | 92.35% | 11.807% |
| 24 | 13.145 | 1841 | 1846 | 1850 | rBV2 | 89515 | 146576 | 3.17% | 0.405% |
| 25 | 13.233 | 1858 | 1861 | 1862 | rBV2 | 75729 | 82013 | 1.77% | 0.227% |
| 26 | 13.257 | 1862 | 1865 | 1868 | rVB | 180921 | 174812 | 3.78% | 0.483% |
| 27 | 13.321 | 1873 | 1876 | 1879 | rBV | 173699 | 156829 | 3.39% | 0.433% |
| 28 | 13.357 | 1879 | 1882 | 1885 | rVB2 | 103416 | 112014 | 2.42% | 0.310% |
| 29 | 13.451 | 1895 | 1898 | 1901 | rBV | 55098 | 55087 | 1.19% | 0.152% |
| 30 | 13.480 | 1901 | 1903 | 1907 | rBV2 | 53535 | 63592 | 1.37% | 0.176% |
| 31 | 13.880 | 1966 | 1971 | 1973 | rBV2 | 55052 | 66154 | 1.43% | 0.183% |
| 32 | 13.904 | 1973 | 1975 | 1977 | rVV | 77032 | 66889 | 1.45% | 0.185% |
| 33 | 13.927 | 1977 | 1979 | 1983 | rVV2 | 74460 | 85299 | 1.84% | 0.236% |
| 34 | 14.004 | 1987 | 1992 | 2005 | rVV5 | 182818 | 554252 | 11.98% | 1.532% |
| 35 | 14.127 | 2005 | 2013 | 2016 | rVV2 | 1138011 | 1411225 | 30.50% | 3.900% |
| 36 | 14.151 | 2016 | 2017 | 2023 | rVB | 250621 | 233540 | 5.05% | 0.645% |

LSC Area Percent Report

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Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

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-BF040622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 14.233 | 2025 | 2031 | 2034 | rBV2 | 109150 | 162022 | 3.50% | 0.448% |
| 38 | 14.515 | 2076 | 2079 | 2082 | rVB | 69325 | 59052 | 1.28% | 0.163% |
| 39 | 14.751 | 2116 | 2119 | 2125 | rVB2 | 55092 | 54806 | 1.18% | 0.151% |
| 40 | 15.015 | 2161 | 2164 | 2168 | rVB4 | 40894 | 52925 | 1.14% | 0.146% |
| 41 | 15.192 | 2190 | 2194 | 2198 | rBV | 229059 | 361284 | 7.81% | 0.998% |
| 42 | 15.263 | 2203 | 2206 | 2211 | rVB4 | 39442 | 51017 | 1.10% | 0.141% |
| 43 | 15.515 | 2245 | 2249 | 2255 | rVB | 114674 | 146063 | 3.16% | 0.404% |
| 44 | 15.580 | 2256 | 2260 | 2266 | rVB | 127875 | 164700 | 3.56% | 0.455% |
| 45 | 15.651 | 2266 | 2272 | 2276 | rBV | 646684 | 870673 | 18.82% | 2.406% |
| 46 | 17.192 | 2528 | 2534 | 2542 | rVB2 | 50334 | 111675 | 2.41% | 0.309% |
| 47 | 17.656 | 2610 | 2613 | 2618 | rVB2 | 30470 | 53902 | 1.17% | 0.149% |

Sum of corrected areas: 36184438

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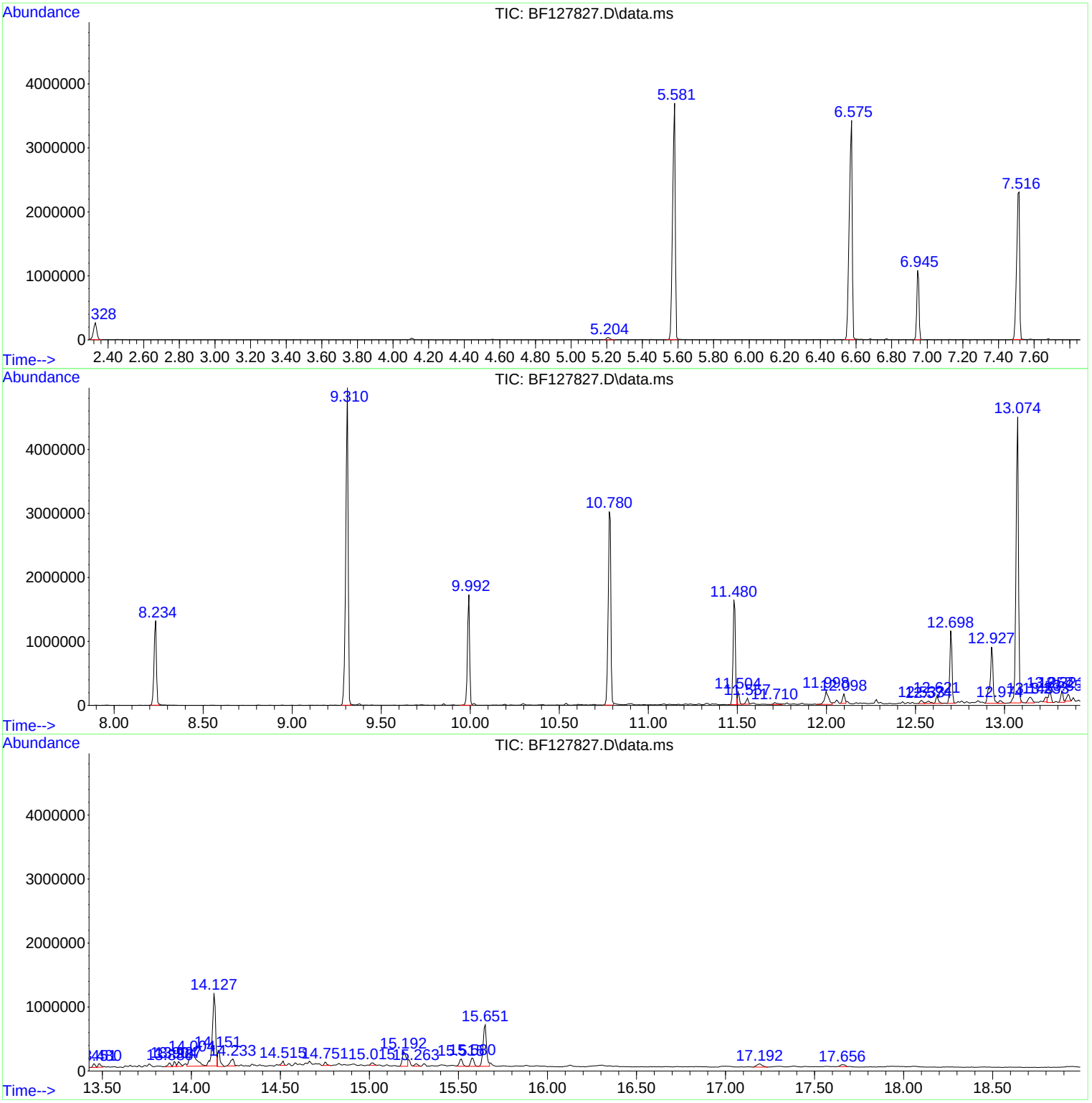
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

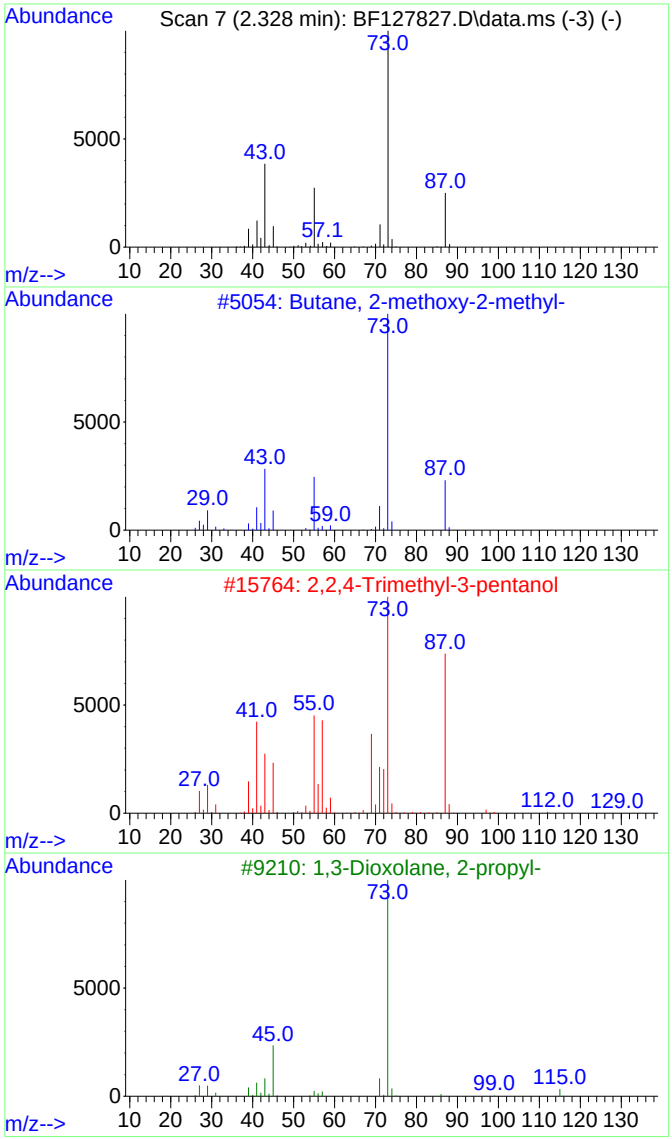
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

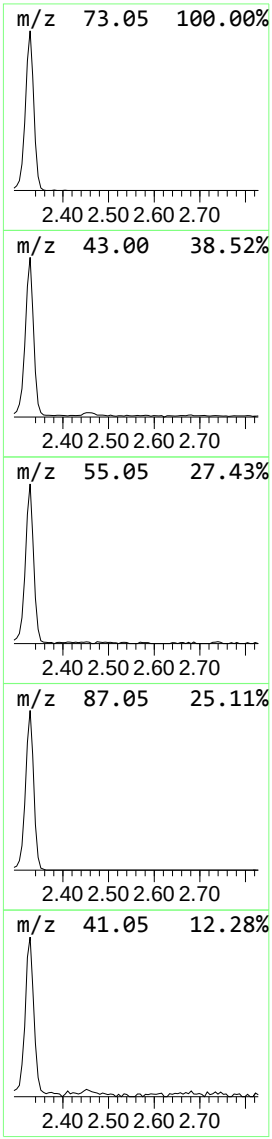
**********

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

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.328 | 6.93 ng | 329646 | 1,4-Dichlorobenzene-d4 | 6.945 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 83 |
| 2 | 2,2,4-Trimethyl-3-pentanol | 130 | C8H18O | 005162-48-1 | 40 |
| 3 | 1,3-Dioxolane, 2-propyl- | 116 | C6H12O2 | 003390-13-4 | 23 |
| 4 | 1,3-Dioxolane, 2-methyl- | 88 | C4H8O2 | 000497-26-7 | 23 |
| 5 | Acetamide, N-ethyl- | 87 | C4H9NO | 000625-50-3 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

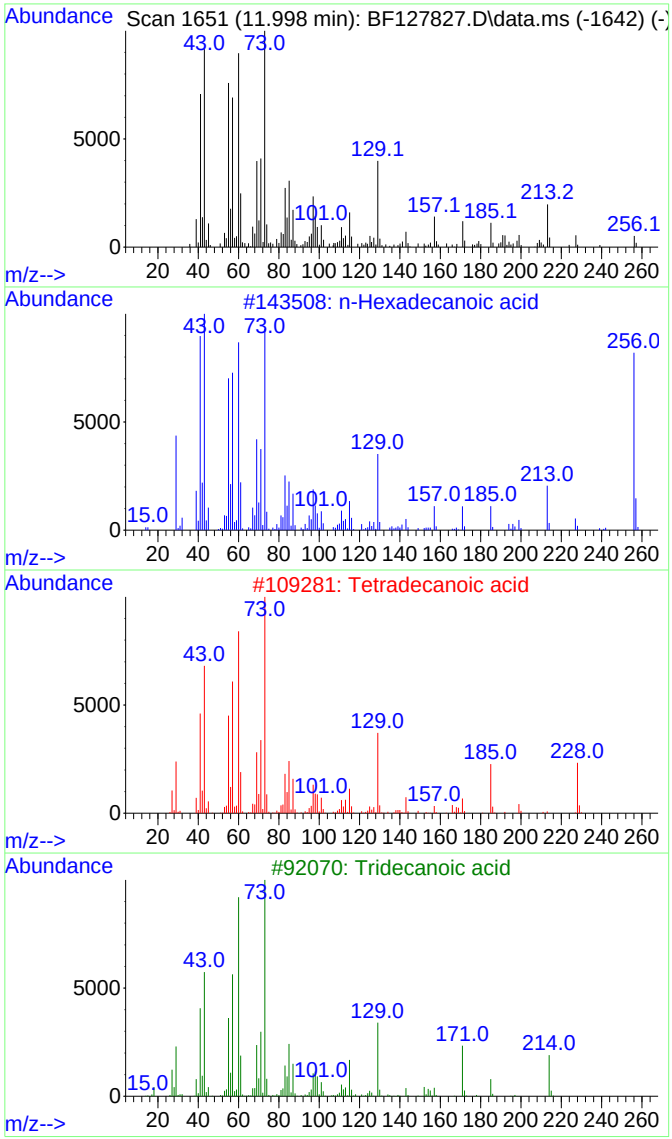
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

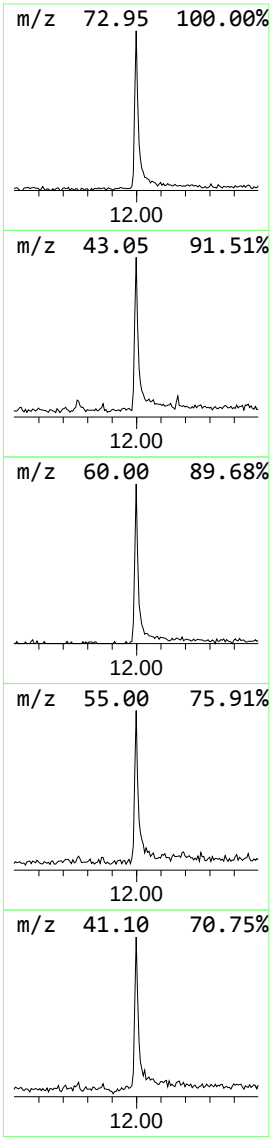
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.998 | 4.17 ng | 315856 | Phenanthrene-d10 | 11.480 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98 |
| 2 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 91 |
| 3 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 87 |
| 4 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 83 |
| 5 | n-Decanoic acid | 172 | C10H20O2 | 000334-48-5 | 62 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

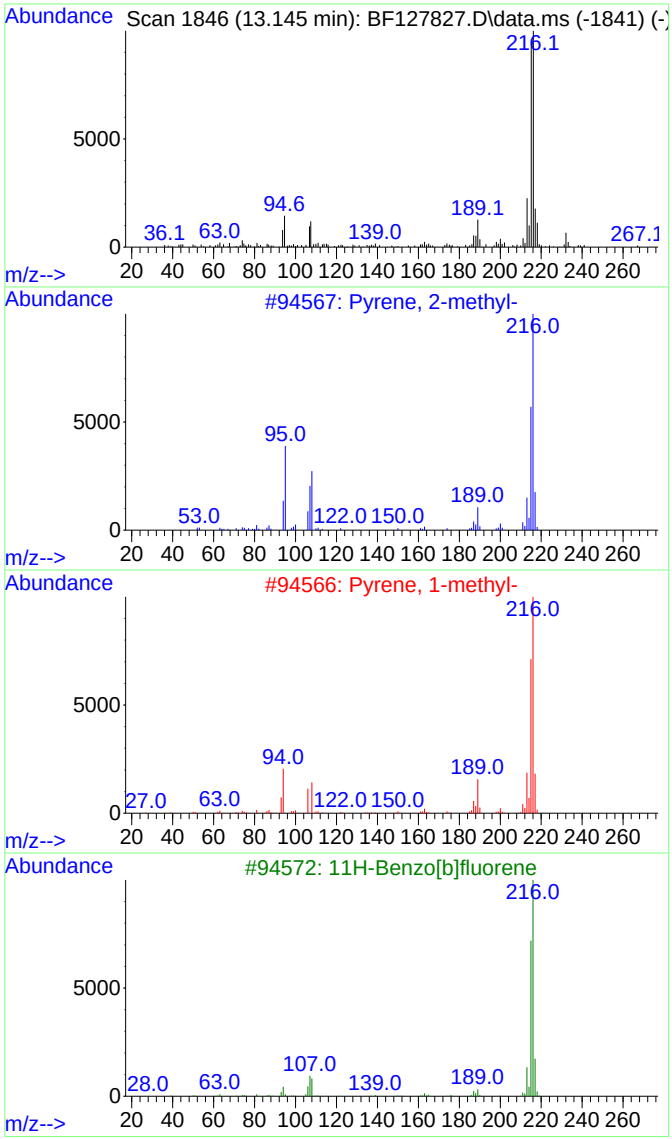
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

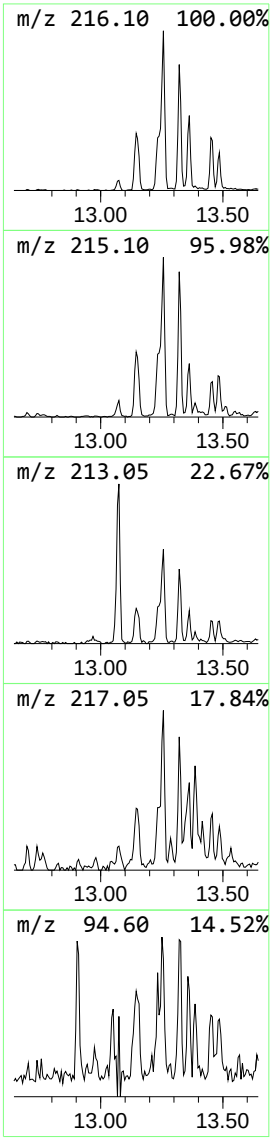
**********

Peak Number 3 Pyrene, 2-methyl- Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.145 | 2.08 ng | 146576 | Chrysene-d12 | 14.127 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 92 |
| 2 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 92 |
| 3 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 87 |
| 4 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 76 |
| 5 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 76 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

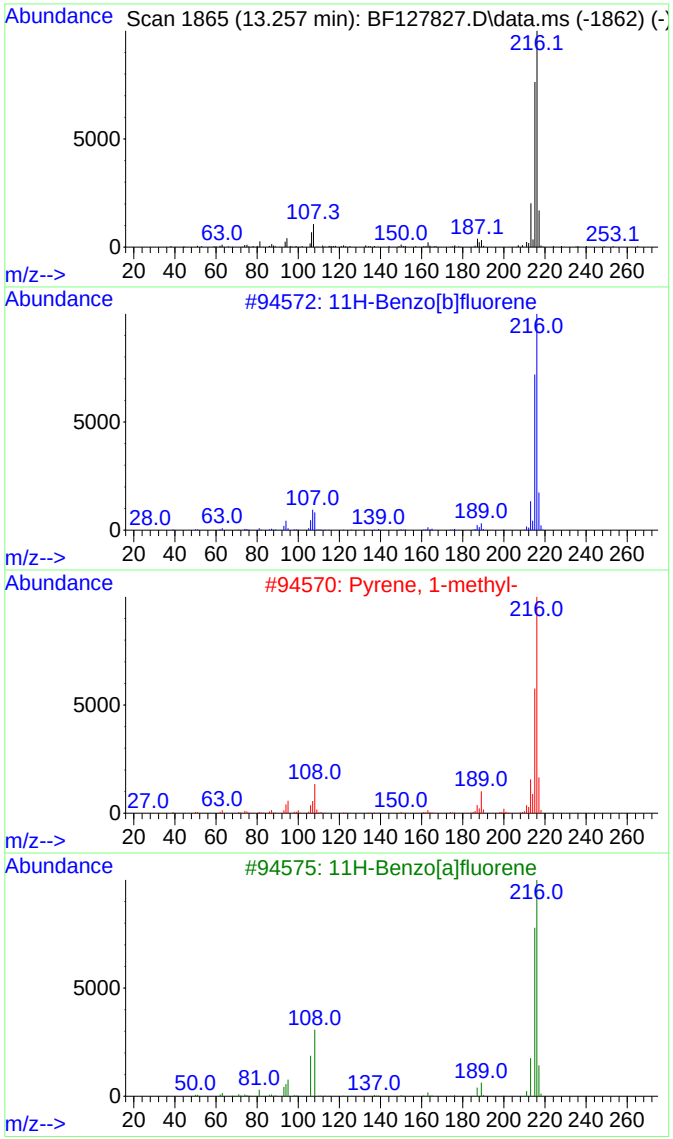
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

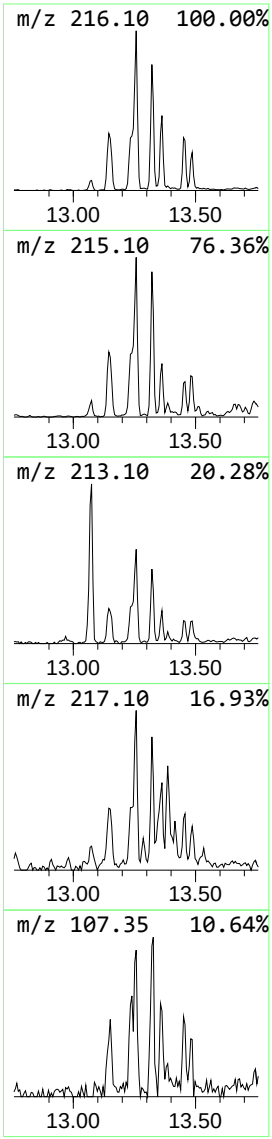
**********

Peak Number 4 11H-Benzo[b]fluorene Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.257 | 2.48 ng | 174812 | Chrysene-d12 | 14.127 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 97 |
| 2 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 91 |
| 3 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 87 |
| 4 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 83 |
| 5 | 1,3,5-Triazine-2,4,6-triamine, N... | 216 | C10H12N6 | 046731-79-7 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

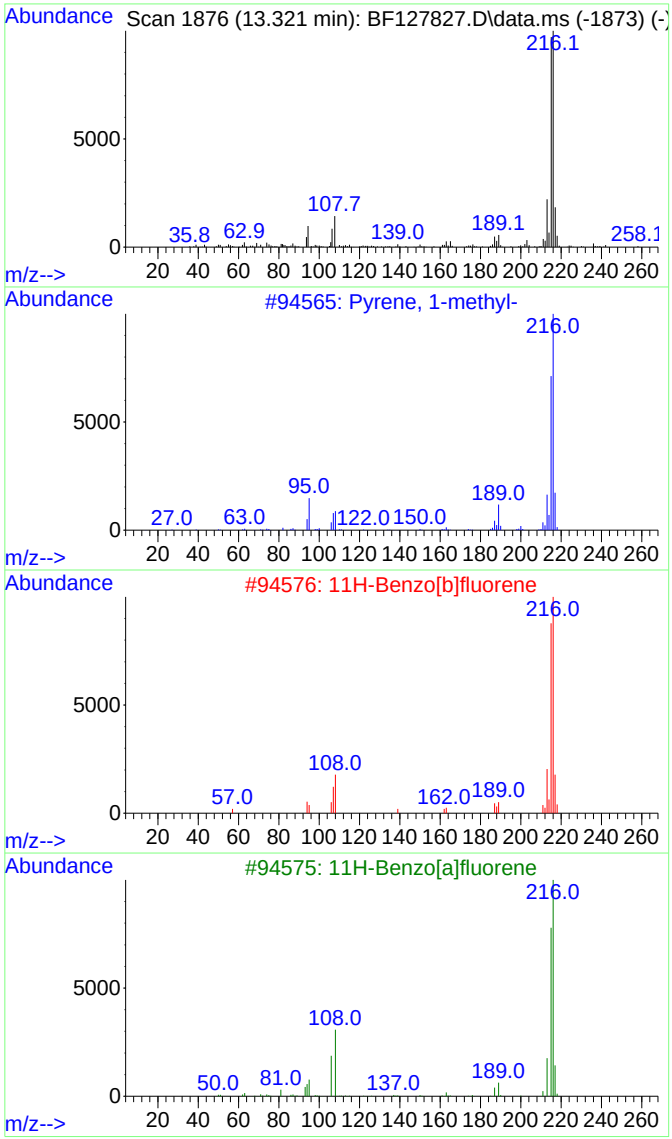
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

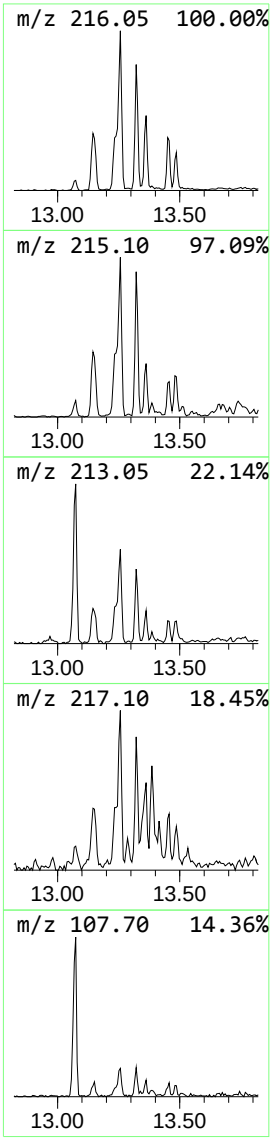
**********

Peak Number 5 Pyrene, 1-methyl- Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.322 | 2.22 ng | 156829 | Chrysene-d12 | 14.127 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 93 |
| 2 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 91 |
| 3 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 87 |
| 4 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 87 |
| 5 | 7H-Benzo[c]fluorene | 216 | C17H12 | 000205-12-9 | 68 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

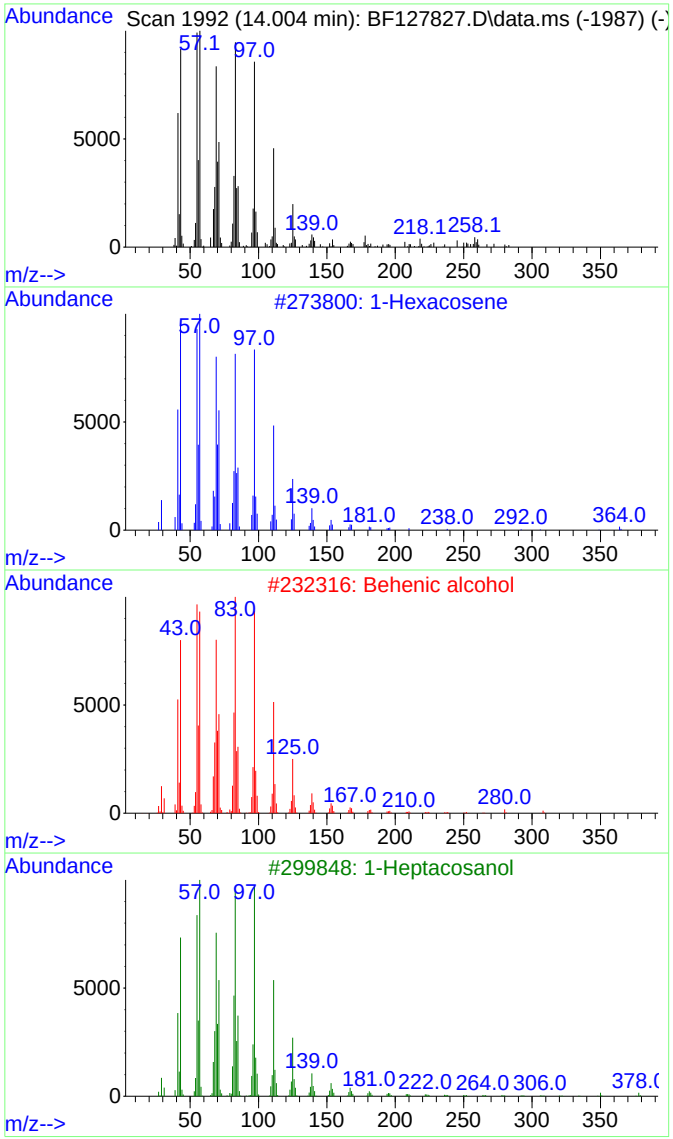
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

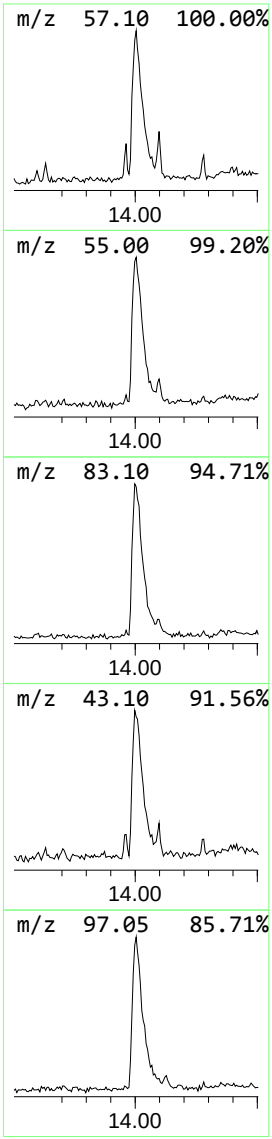
**********

Peak Number 6 1-Hexacosene Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.004 | 7.85 ng | 554252 | Chrysene-d12 | 14.127 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Hexacosene | 364 | C26H52 | 018835-33-1 | 94 |
| 2 | Behenic alcohol | 326 | C22H46O | 000661-19-8 | 91 |
| 3 | 1-Heptacosanol | 396 | C27H56O | 002004-39-9 | 91 |
| 4 | 1-Tetracosene | 336 | C24H48 | 010192-32-2 | 91 |
| 5 | Trifluoroacetoxy hexadecane | 338 | C18H33F3O2 | 006222-03-3 | 90 |



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Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

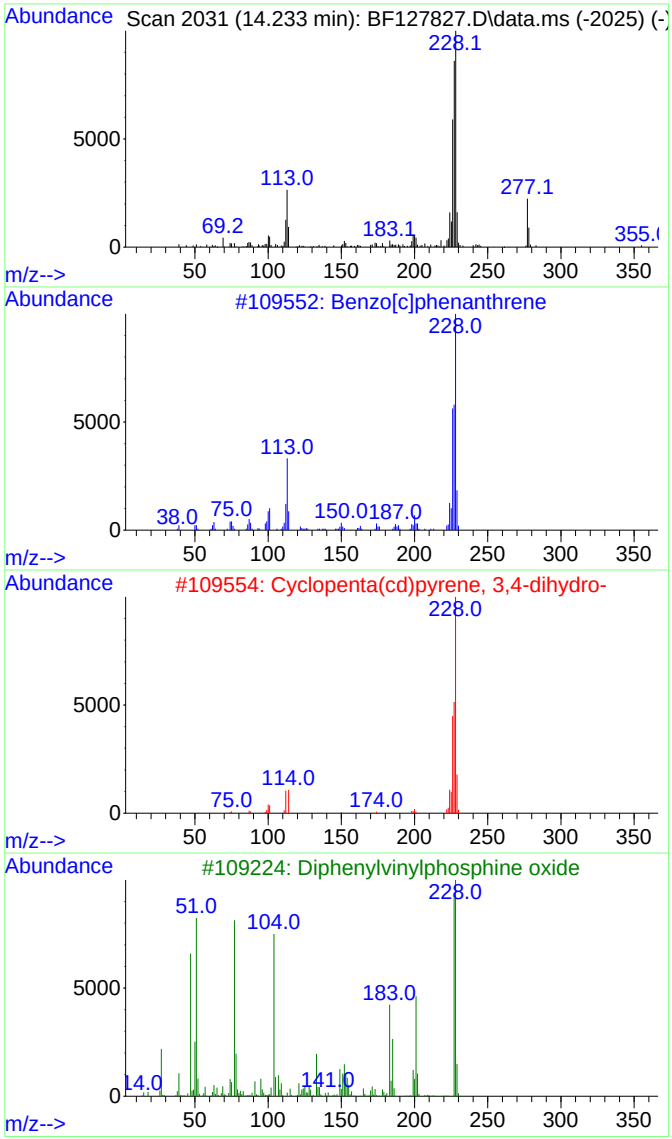
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

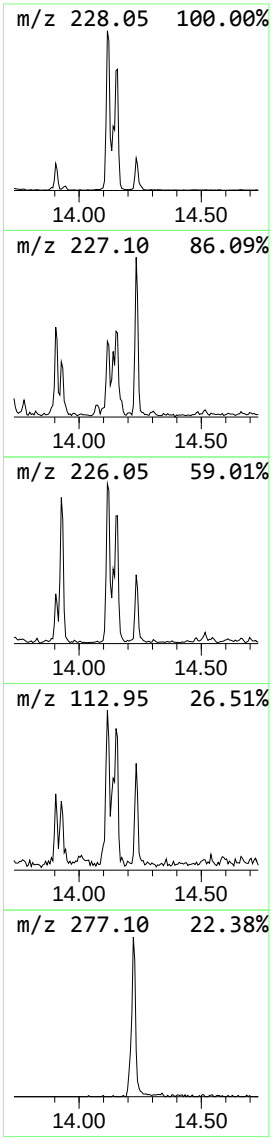
**********

Peak Number 7 Benzo[c]phenanthrene Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.233 | 2.30 ng | 162022 | Chrysene-d12 | 14.127 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[c]phenanthrene | 228 | C18H12 | 000195-19-7 | 55 |
| 2 | Cyclopenta(cd)pyrene, 3,4-dihydro- | 228 | C18H12 | 025732-74-5 | 50 |
| 3 | Diphenylvinylphosphine oxide | 228 | C14H13OP | 002096-78-8 | 43 |
| 4 | Benz[a]anthracene | 228 | C18H12 | 000056-55-3 | 38 |
| 5 | Chrysene | 228 | C18H12 | 000218-01-9 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

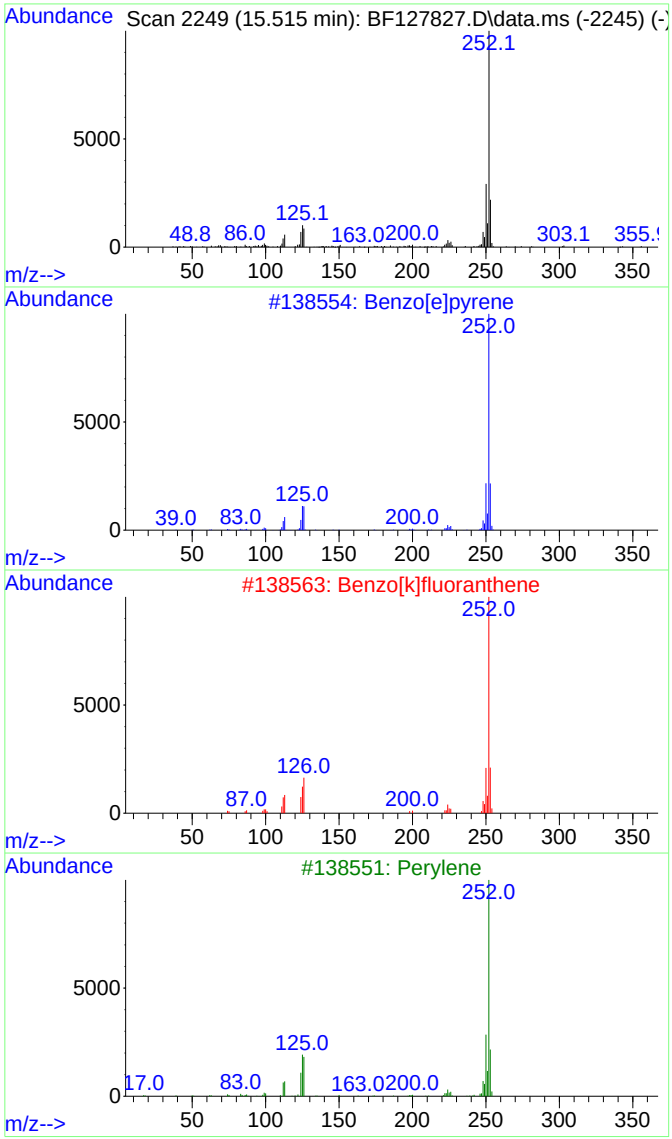
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

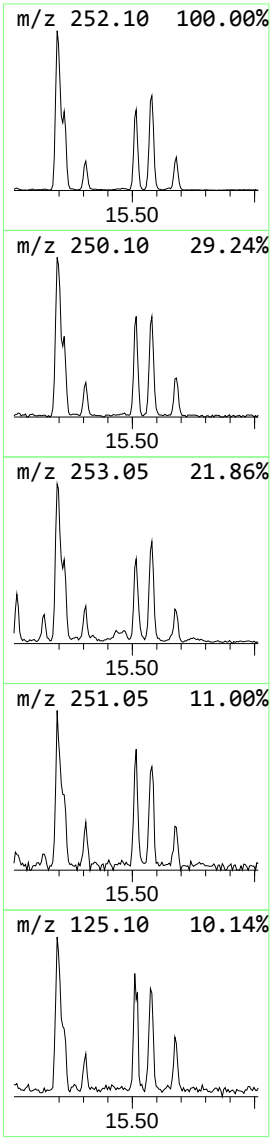
**********

Peak Number 8 Benzo[e]pyrene Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.515 | 3.36 ng | 146063 | Perylene-d12 | 15.651 |

| 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 | Perylene | 252 | C20H12 | 000198-55-0 | 97 |
| 4 | Benzo[j]fluoranthene | 252 | C20H12 | 000205-82-3 | 94 |
| 5 | Benzo[b]fluoranthene | 252 | C20H12 | 000205-99-2 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

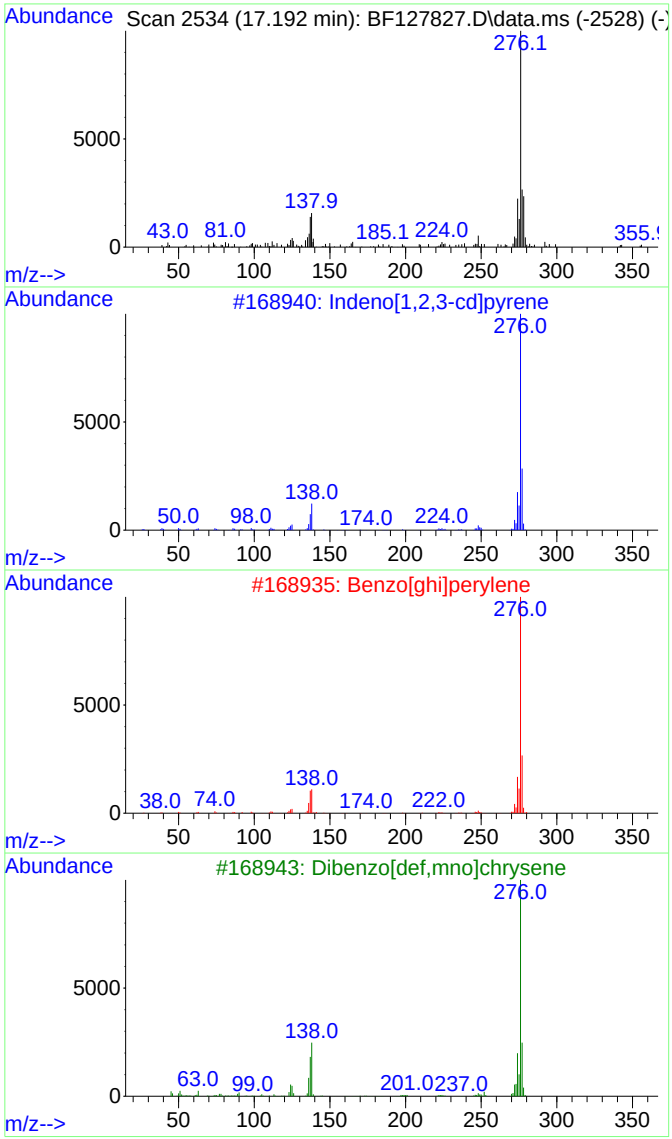
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

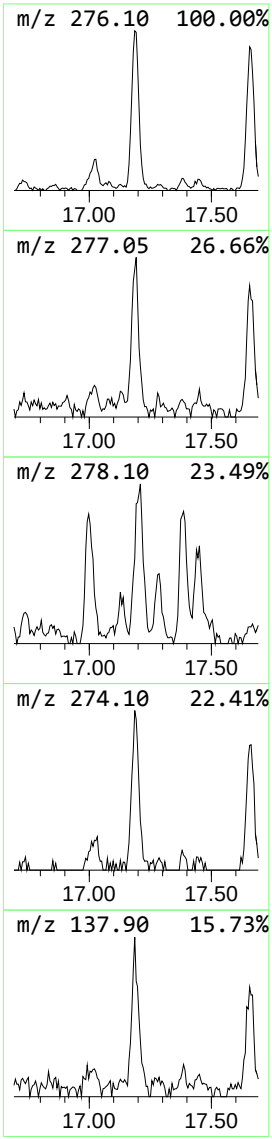
**********

Peak Number 9 Dibenzo[def,mno]chrysene Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.192 | 2.57 ng | 111675 | Perylene-d12 | 15.651 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Indeno[1,2,3-cd]pyrene | 276 | C22H12 | 000193-39-5 | 95 |
| 2 | Benzo[ghi]perylene | 276 | C22H12 | 000191-24-2 | 89 |
| 3 | Dibenzo[def,mno]chrysene | 276 | C22H12 | 000191-26-4 | 70 |
| 4 | 12H-Benzothieno[2,3-d]pyrido[1,2... | 276 | C14H16N2S2 | 102254-63-7 | 50 |
| 5 | Benzofuro[3,2-d]pyrimidine, 4-me... | 276 | C17H12N2O2 | 312265-36-4 | 47 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
Data File : BF127827.D
Acq On : 18 Apr 2022 16:08
Operator : CG\JU
Sample : N2442-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
124030

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Butane, 2-metho... | 2.328 | 6.9 | ng | 329646 | 1 | 6.945 | 950904 | 20.0 |
| n-Hexadecanoic ... | 11.998 | 4.2 | ng | 315856 | 4 | 11.480 | 1516250 | 20.0 |
| Pyrene, 2-methyl- | 13.145 | 2.1 | ng | 146576 | 5 | 14.127 | 1411230 | 20.0 |
| 11H-Benzo[b]flu... | 13.257 | 2.5 | ng | 174812 | 5 | 14.127 | 1411230 | 20.0 |
| Pyrene, 1-methyl- | 13.322 | 2.2 | ng | 156829 | 5 | 14.127 | 1411230 | 20.0 |
| 1-Hexacosene | 14.004 | 7.8 | ng | 554252 | 5 | 14.127 | 1411230 | 20.0 |
| Benzo[c]phenant... | 14.233 | 2.3 | ng | 162022 | 5 | 14.127 | 1411230 | 20.0 |
| Benzo[e]pyrene | 15.515 | 3.4 | ng | 146063 | 6 | 15.651 | 870673 | 20.0 |
| Dibenzo[def,mno... | 17.192 | 2.6 | ng | 111675 | 6 | 15.651 | 870673 | 20.0 |