LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

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-BF072522.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF129445.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 5.510 | 545 | 548 | 554 | rVB | 1377892 | 1277018 | 1.80% | 0.200% |
| 2 | 5.846 | 600 | 605 | 608 | rBV | 2398021 | 2732596 | 3.85% | 0.428% |
| 3 | 6.522 | 717 | 720 | 724 | rBV | 728317 | 732959 | 1.03% | 0.115% |
| 4 | 6.634 | 734 | 739 | 742 | rVB | 581564 | 957924 | 1.35% | 0.150% |
| 5 | 6.881 | 778 | 781 | 784 | rBV | 1037366 | 892328 | 1.26% | 0.140% |
| 6 | 6.957 | 788 | 794 | 797 | rBV | 6688095 | 9074086 | 12.78% | 1.422% |
| 7 | 7.451 | 873 | 878 | 880 | rBV | 2820511 | 2737719 | 3.86% | 0.429% |
| 8 | 7.928 | 891 | 959 | 961 | rBV | 6880730 | 71004871 | 100.00% | 11.125% |
| 9 | 8.093 | 975 | 987 | 988 | rBV2 | 1585030 | 3394961 | 4.78% | 0.532% |
| 10 | 8.140 | 988 | 995 | 997 | rBV | 1685332 | 3012441 | 4.24% | 0.472% |
| 11 | 8.169 | 997 | 1000 | 1002 | rVV | 1284845 | 1431936 | 2.02% | 0.224% |
| 12 | 8.751 | 1095 | 1099 | 1102 | rVB | 3636582 | 2906049 | 4.09% | 0.455% |
| 13 | 9.187 | 1156 | 1173 | 1175 | rBV | 3082643 | 8333480 | 11.74% | 1.306% |
| 14 | 9.245 | 1179 | 1183 | 1186 | rBV | 4849820 | 4082923 | 5.75% | 0.640% |
| 15 | 9.316 | 1192 | 1195 | 1198 | rVB | 2618732 | 1944257 | 2.74% | 0.305% |
| 16 | 9.528 | 1229 | 1231 | 1237 | rVV | 1256791 | 1529466 | 2.15% | 0.240% |
| 17 | 9.739 | 1264 | 1267 | 1273 | rBV2 | 4245482 | 4828259 | 6.80% | 0.757% |
| 18 | 9.845 | 1282 | 1285 | 1289 | rVV | 2489790 | 2247185 | 3.16% | 0.352% |
| 19 | 9.928 | 1296 | 1299 | 1302 | rVV2 | 2191045 | 2451067 | 3.45% | 0.384% |
| 20 | 9.963 | 1302 | 1305 | 1309 | rVB | 2954797 | 2813708 | 3.96% | 0.441% |
| 21 | 10.045 | 1317 | 1319 | 1327 | rVB | 1211750 | 1561249 | 2.20% | 0.245% |
| 22 | 10.234 | 1333 | 1351 | 1352 | rBV | 6268554 | 18390417 | 25.90% | 2.881% |
| 23 | 10.251 | 1352 | 1354 | 1363 | rVB | 4938456 | 5918088 | 8.33% | 0.927% |
| 24 | 10.551 | 1402 | 1405 | 1410 | rVB | 849755 | 1075754 | 1.52% | 0.169% |
| 25 | 10.610 | 1412 | 1415 | 1418 | rVB | 1069009 | 809642 | 1.14% | 0.127% |
| 26 | 10.739 | 1431 | 1437 | 1443 | rBV2 | 4690424 | 6774328 | 9.54% | 1.061% |
| 27 | 10.816 | 1447 | 1450 | 1453 | rBV | 911342 | 740032 | 1.04% | 0.116% |
| 28 | 10.951 | 1464 | 1473 | 1475 | rBV2 | 5865155 | 8225575 | 11.58% | 1.289% |
| 29 | 10.981 | 1475 | 1478 | 1481 | rVV | 10801600 | 14132989 | 19.90% | 2.214% |
| 30 | 11.004 | 1481 | 1482 | 1491 | rVB | 7757012 | 8790490 | 12.38% | 1.377% |
| 31 | 11.104 | 1495 | 1499 | 1503 | rBV3 | 601167 | 1166734 | 1.64% | 0.183% |
| 32 | 11.198 | 1513 | 1515 | 1520 | rBV | 4278463 | 4654364 | 6.55% | 0.729% |
| 33 | 11.251 | 1523 | 1524 | 1526 | rVV | 993488 | 922172 | 1.30% | 0.144% |
| 34 | 11.286 | 1526 | 1530 | 1539 | rVB | 8920578 | 11974506 | 16.86% | 1.876% |
| 35 | 11.363 | 1540 | 1543 | 1546 | rBV | 2589679 | 2099619 | 2.96% | 0.329% |
| 36 | 11.422 | 1549 | 1553 | 1556 | rBV | 1125080 | 1042737 | 1.47% | 0.163% |

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Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

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-BF072522.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 11.539 | 1568 | 1573 | 1577 | rBV3 | 778679 | 1501927 | 2.12% | 0.235% |
| 38 | 11.633 | 1586 | 1589 | 1593 | rBV2 | 1523288 | 2137028 | 3.01% | 0.335% |
| 39 | 11.669 | 1593 | 1595 | 1598 | rVB2 | 897886 | 800849 | 1.13% | 0.125% |
| 40 | 11.798 | 1612 | 1617 | 1620 | rBV3 | 1110204 | 1861315 | 2.62% | 0.292% |
| 41 | 11.857 | 1623 | 1627 | 1629 | rBV2 | 795495 | 1033295 | 1.46% | 0.162% |
| 42 | 11.945 | 1638 | 1642 | 1644 | rBV2 | 2120901 | 3127938 | 4.41% | 0.490% |
| 43 | 12.039 | 1656 | 1658 | 1661 | rVB | 1582259 | 1287316 | 1.81% | 0.202% |
| 44 | 12.075 | 1661 | 1664 | 1666 | rBV2 | 1416055 | 1606475 | 2.26% | 0.252% |
| 45 | 12.128 | 1668 | 1673 | 1676 | rBV2 | 1970439 | 3580128 | 5.04% | 0.561% |
| 46 | 12.216 | 1680 | 1688 | 1690 | rBV | 12229524 | 22268051 | 31.36% | 3.489% |
| 47 | 12.433 | 1717 | 1725 | 1729 | rBV6 | 6549634 | 18139601 | 25.55% | 2.842% |
| 48 | 12.580 | 1748 | 1750 | 1752 | rVB2 | 7982605 | 6965078 | 9.81% | 1.091% |
| 49 | 12.610 | 1752 | 1755 | 1758 | rBV2 | 6200273 | 10359622 | 14.59% | 1.623% |
| 50 | 12.722 | 1771 | 1774 | 1775 | rBV | 1865173 | 1736873 | 2.45% | 0.272% |
| 51 | 12.810 | 1775 | 1789 | 1794 | rVV7 | 13358696 | 65356989 | 92.05% | 10.240% |
| 52 | 12.863 | 1796 | 1798 | 1803 | rVV2 | 8635018 | 11819552 | 16.65% | 1.852% |
| 53 | 12.904 | 1803 | 1805 | 1808 | rVB3 | 9568689 | 9139764 | 12.87% | 1.432% |
| 54 | 12.939 | 1808 | 1811 | 1818 | rVB2 | 3966225 | 6179455 | 8.70% | 0.968% |
| 55 | 13.010 | 1818 | 1823 | 1824 | rBV | 15190505 | 22490837 | 31.68% | 3.524% |
| 56 | 13.045 | 1827 | 1829 | 1832 | rVV2 | 1971651 | 1608939 | 2.27% | 0.252% |
| 57 | 13.080 | 1832 | 1835 | 1838 | rVB2 | 4252991 | 4884659 | 6.88% | 0.765% |
| 58 | 13.110 | 1838 | 1840 | 1843 | rVB3 | 5406973 | 4211165 | 5.93% | 0.660% |
| 59 | 13.133 | 1843 | 1844 | 1846 | rVB | 3908992 | 2280525 | 3.21% | 0.357% |
| 60 | 13.180 | 1846 | 1852 | 1853 | rBV3 | 5308488 | 9020319 | 12.70% | 1.413% |
| 61 | 13.233 | 1859 | 1861 | 1865 | rVV3 | 2885584 | 2830909 | 3.99% | 0.444% |
| 62 | 13.269 | 1865 | 1867 | 1872 | rVB2 | 4271758 | 4306986 | 6.07% | 0.675% |
| 63 | 13.369 | 1876 | 1884 | 1886 | rBV | 14939646 | 31060143 | 43.74% | 4.867% |
| 64 | 13.392 | 1886 | 1888 | 1891 | rVB2 | 7767455 | 6166336 | 8.68% | 0.966% |
| 65 | 13.439 | 1892 | 1896 | 1901 | rBV2 | 5894454 | 8589432 | 12.10% | 1.346% |
| 66 | 13.516 | 1907 | 1909 | 1912 | rVB | 2308305 | 1853521 | 2.61% | 0.290% |
| 67 | 13.563 | 1915 | 1917 | 1919 | rBV2 | 994270 | 931416 | 1.31% | 0.146% |
| 68 | 13.663 | 1931 | 1934 | 1935 | rBV | 3184566 | 2867833 | 4.04% | 0.449% |
| 69 | 13.710 | 1935 | 1942 | 1944 | rVV | 14371613 | 28089930 | 39.56% | 4.401% |
| 70 | 13.780 | 1951 | 1954 | 1957 | rBV | 3636533 | 3652556 | 5.14% | 0.572% |
| 71 | 13.916 | 1973 | 1977 | 1985 | rVB | 3678326 | 5098657 | 7.18% | 0.799% |
| 72 | 14.045 | 1991 | 1999 | 2002 | rBV | 14513095 | 26795115 | 37.74% | 4.198% |
| 73 | 14.280 | 2035 | 2039 | 2042 | rBV | 2745235 | 2664029 | 3.75% | 0.417% |
| 74 | 14.527 | 2077 | 2081 | 2084 | rVB2 | 2531936 | 2427262 | 3.42% | 0.380% |
| 75 | 14.580 | 2087 | 2090 | 2096 | rVB | 736347 | 952187 | 1.34% | 0.149% |
| 76 | 14.674 | 2103 | 2106 | 2113 | rVB2 | 608445 | 933603 | 1.31% | 0.146% |

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Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

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-BF072522.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 14.733 | 2113 | 2116 | 2119 | rVB | 803716 | 757690 | 1.07% | 0.119% |
| 78 | 14.957 | 2151 | 2154 | 2159 | rVB | 1078036 | 1231984 | 1.74% | 0.193% |
| 79 | 15.180 | 2188 | 2192 | 2195 | rVV2 | 647162 | 930313 | 1.31% | 0.146% |
| 80 | 15.216 | 2195 | 2198 | 2208 | rVB2 | 894161 | 1637122 | 2.31% | 0.257% |
| 81 | 15.433 | 2231 | 2235 | 2240 | rVB | 1827753 | 2340354 | 3.30% | 0.367% |
| 82 | 15.568 | 2254 | 2258 | 2268 | rVB | 723418 | 1082124 | 1.52% | 0.170% |
| 83 | 16.510 | 2402 | 2418 | 2422 | rBV | 12212155 | 39889413 | 56.18% | 6.250% |
| 84 | 17.792 | 2620 | 2636 | 2657 | rVB2 | 7986062 | 31081169 | 43.77% | 4.870% |

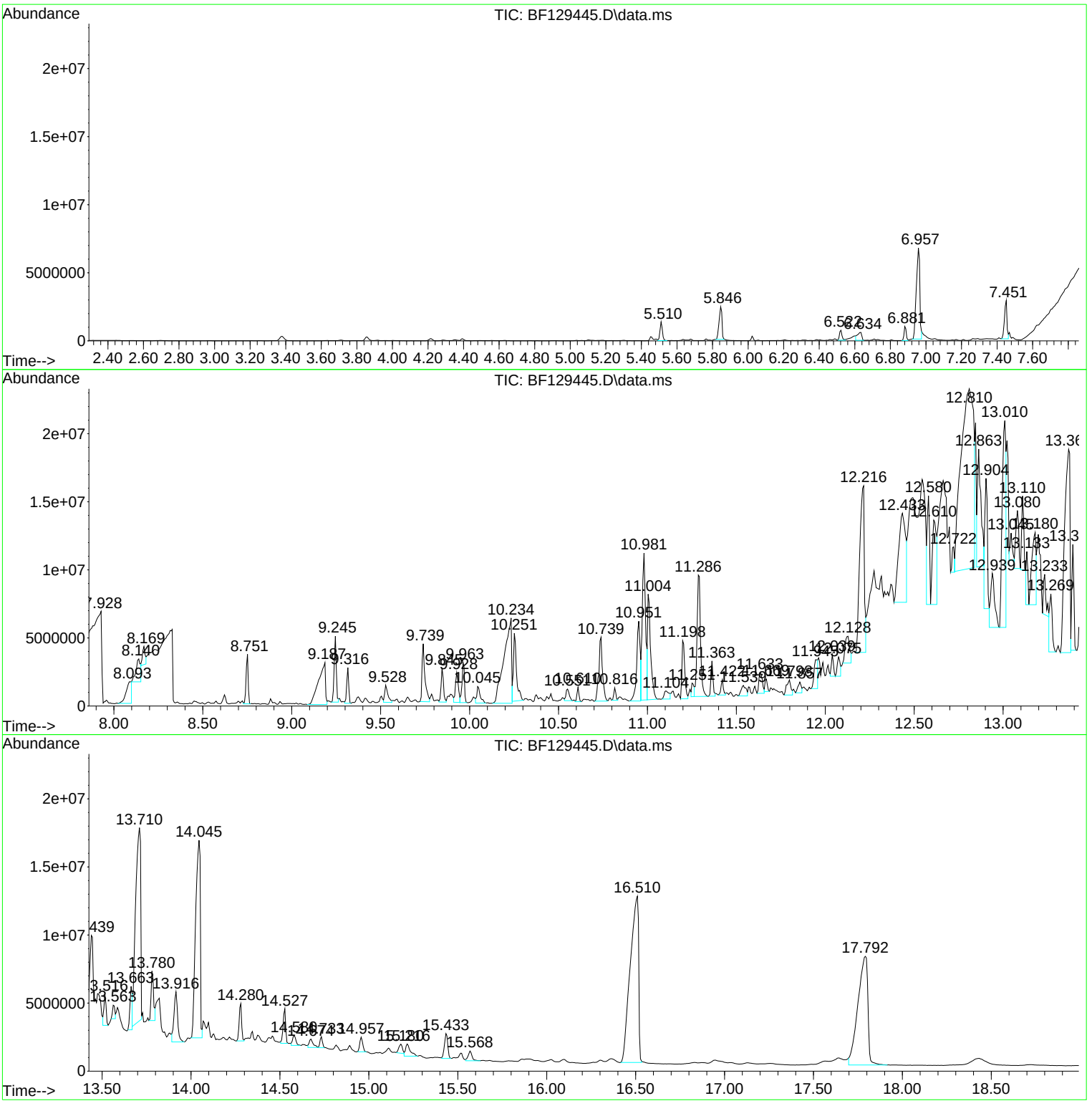
Sum of corrected areas: 638231713

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.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\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

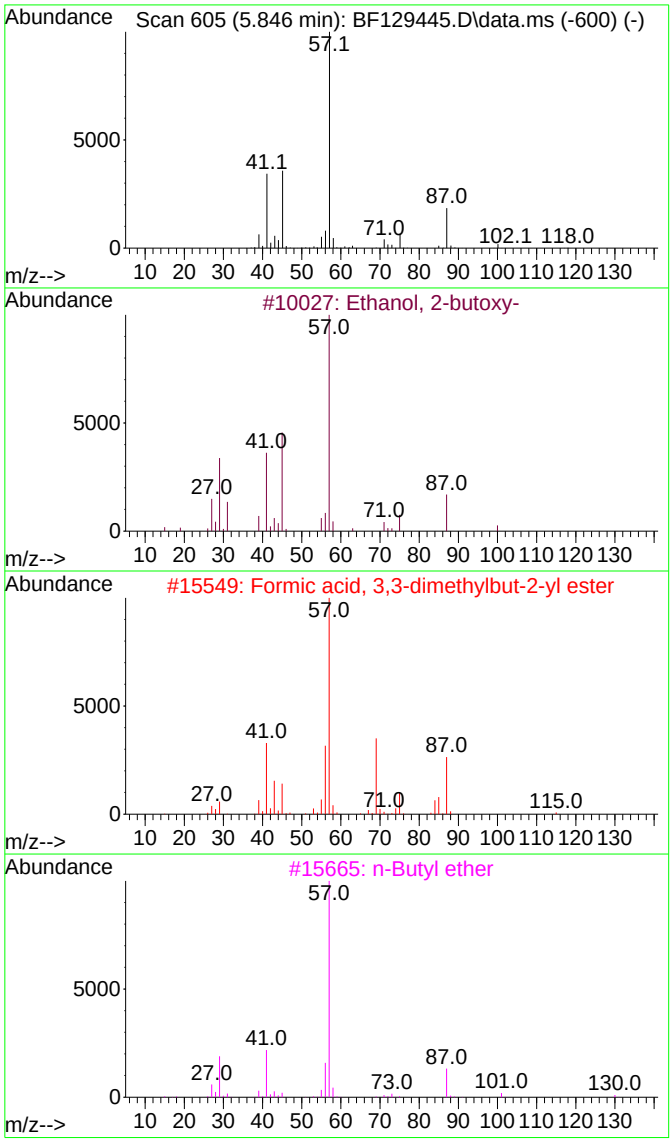
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

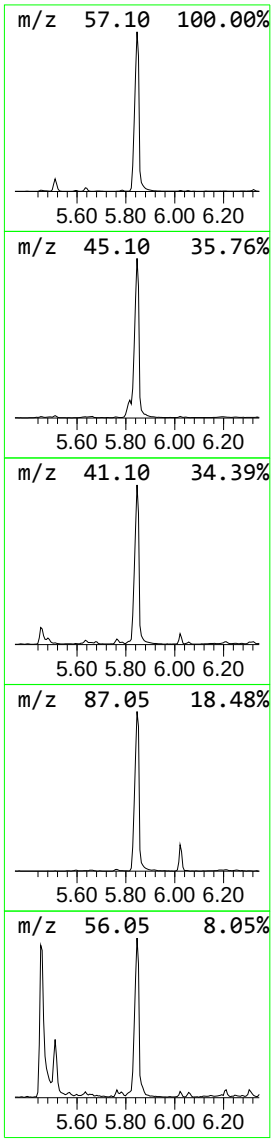
**********

Peak Number 1 Ethanol, 2-butoxy- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.846 | 61.25 ng | 2732600 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Ethanol, 2-butoxy- | 118 | C6H14O2 | 000111-76-2 | 78 |
| 2 | Formic acid, 3,3-dimethylbut-2-y... | 130 | C7H14O2 | 1000368-20-5 | 33 |
| 3 | n-Butyl ether | 130 | C8H18O | 000142-96-1 | 32 |
| 4 | Propanoic acid, 2-methylpropyl e... | 130 | C7H14O2 | 000540-42-1 | 25 |
| 5 | 1,3-Propanediol, 2-methyl-, dipr... | 202 | C10H18O4 | 054932-83-1 | 23 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

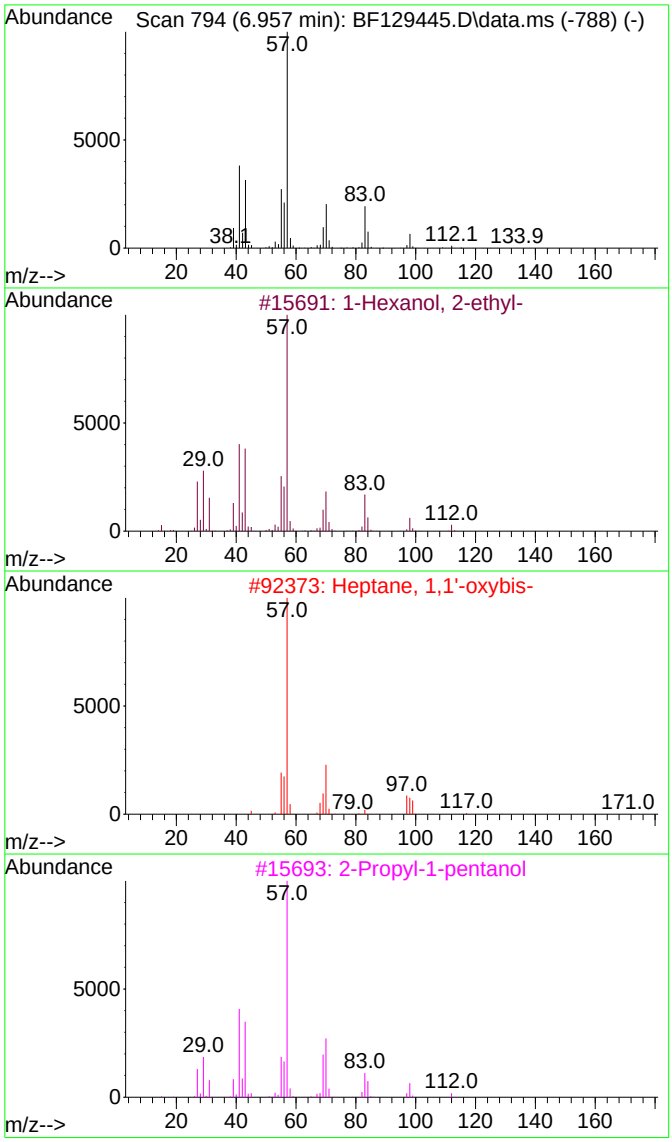
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

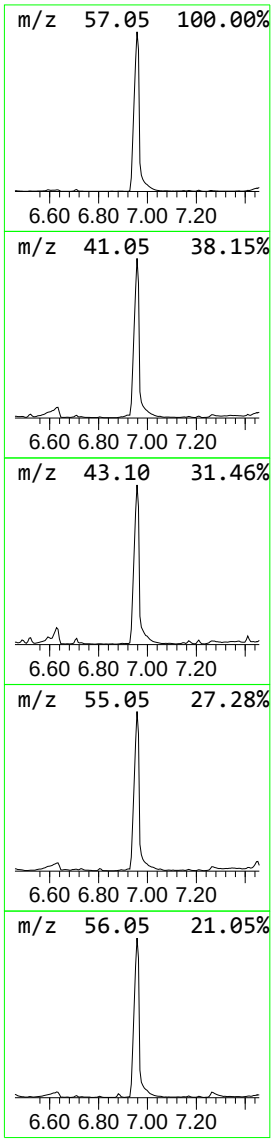
**********

Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.957 | 203.38 ng | 9074090 | 1,4-Dichlorobenzene-d4 | 6.881 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Hexanol, 2-ethyl- | 130 | C8H18O | 000104-76-7 | 78 |
| 2 | Heptane, 1,1'-oxybis- | 214 | C14H30O | 000629-64-1 | 59 |
| 3 | 2-Propyl-1-pentanol | 130 | C8H18O | 058175-57-8 | 47 |
| 4 | Oxalic acid, isobutyl octyl ester | 258 | C14H26O4 | 1000309-37-3 | 43 |
| 5 | Carbonic acid, heptyl vinyl ester | 186 | C10H18O3 | 1010383-25-4 | 43 |



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Acq On : 29 Jul 2022 16:58
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Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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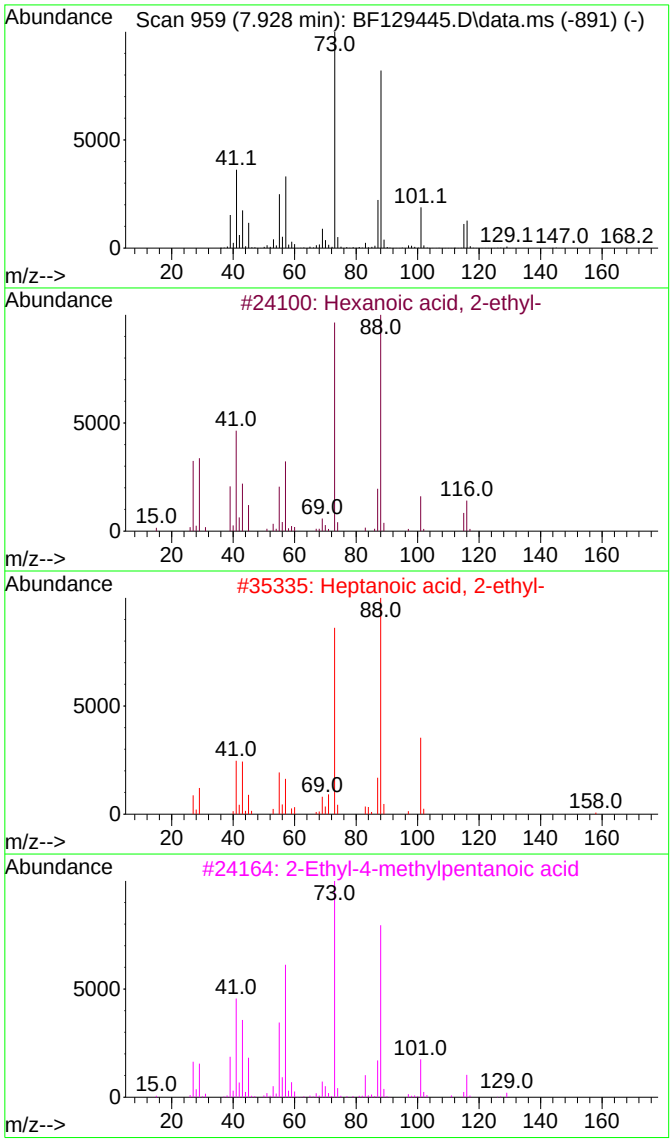
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

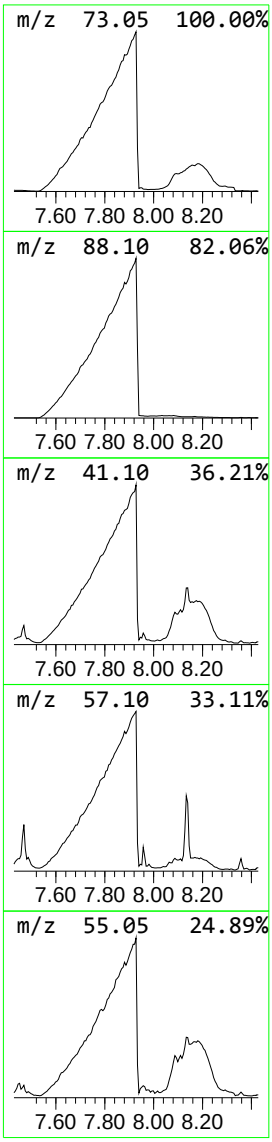
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 3 Hexanoic acid, 2-ethyl- Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.928 | 991.73 ng | 71004900 | Naphthalene-d8 | 8.169 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexanoic acid, 2-ethyl- | 144 | C8H16O2 | 000149-57-5 | 90 |
| 2 | Heptanoic acid, 2-ethyl- | 158 | C9H18O2 | 003274-29-1 | 59 |
| 3 | 2-Ethyl-4-methylpentanoic acid | 144 | C8H16O2 | 000108-81-6 | 56 |
| 4 | Butanoic acid, 2-ethyl-, 1,2,3-p... | 386 | C21H38O6 | 056554-54-2 | 42 |
| 5 | Octanoic acid, 2-methyl-, methyl... | 172 | C10H20O2 | 002177-86-8 | 40 |



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Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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BNA_F
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

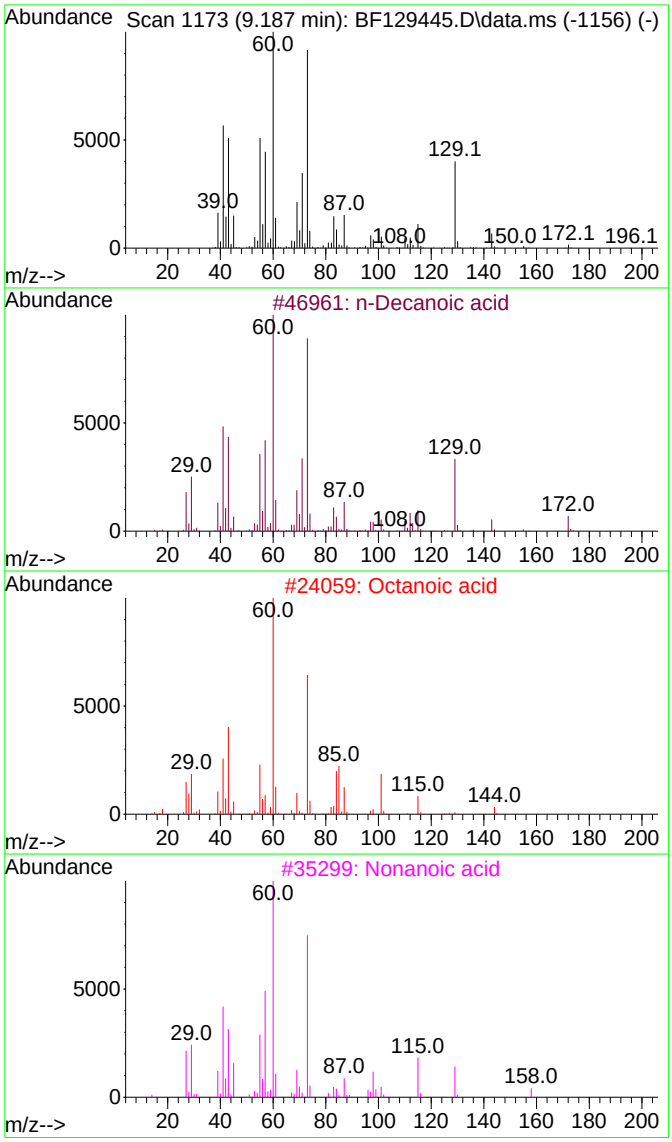
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

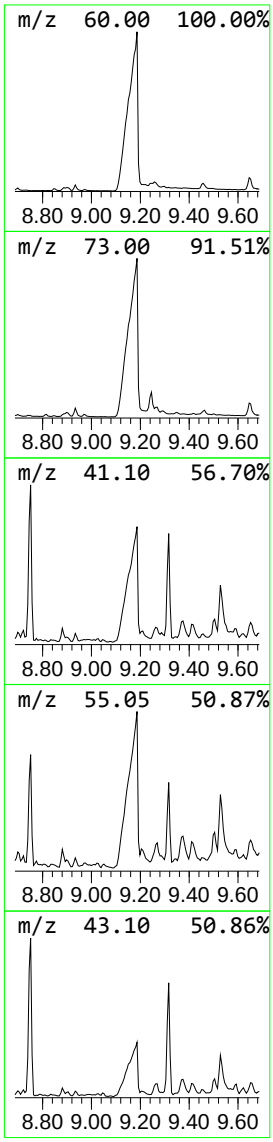
**********

Peak Number 4 n-Decanoic acid Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.187 | 68.00 ng | 8333480 | Acenaphthene-d10 | 9.928 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Decanoic acid | 172 | C10H20O2 | 000334-48-5 | 90 |
| 2 | Octanoic acid | 144 | C8H16O2 | 000124-07-2 | 50 |
| 3 | Nonanoic acid | 158 | C9H18O2 | 000112-05-0 | 50 |
| 4 | Decanoic acid, silver(1+) salt | 278 | C10H19AgO2 | 013126-67-5 | 50 |
| 5 | Propanedioic acid, propyl- | 146 | C6H10O4 | 000616-62-6 | 47 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

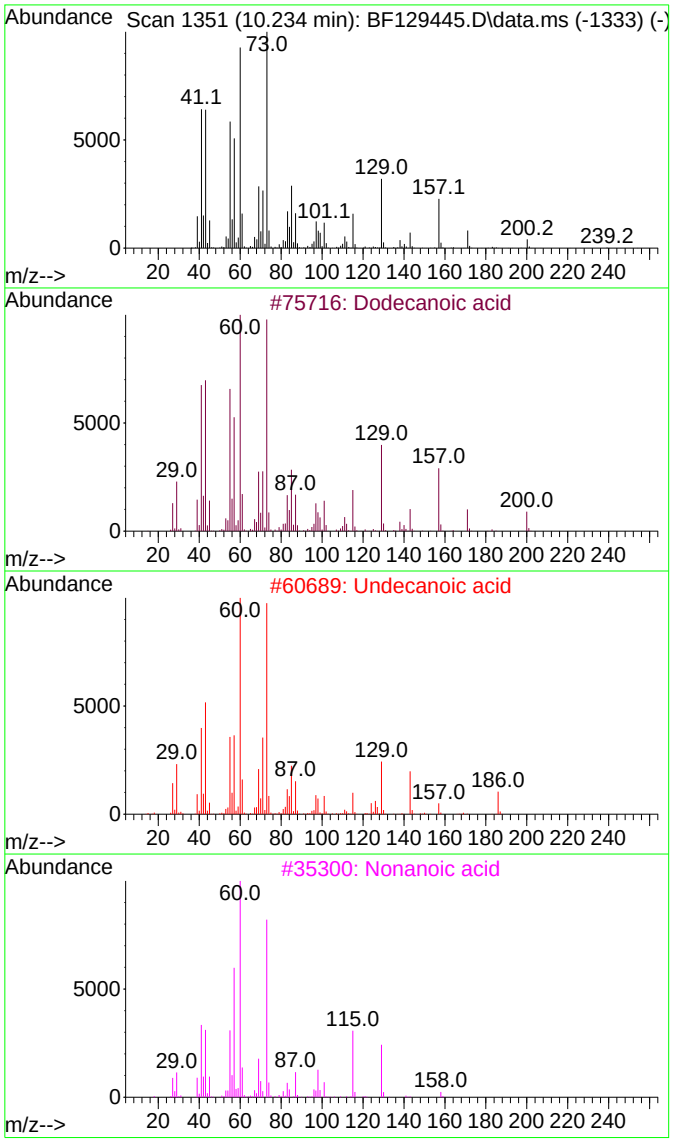
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

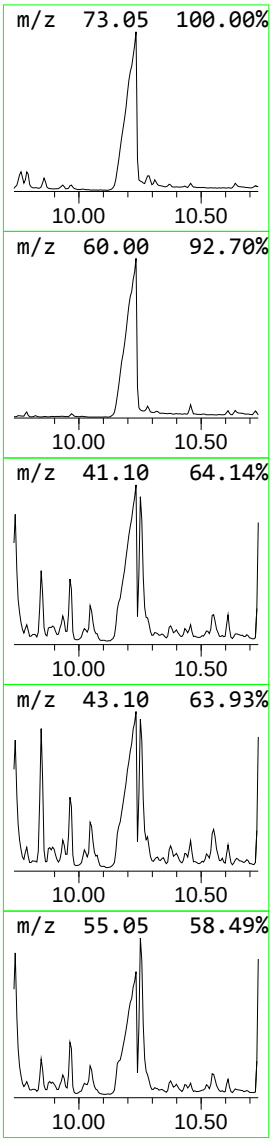
**********

Peak Number 5 Dodecanoic acid Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.233 | 150.06 ng | 18390400 | Acenaphthene-d10 | 9.928 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dodecanoic acid | 200 | C12H24O2 | 000143-07-7 | 97 |
| 2 | Undecanoic acid | 186 | C11H22O2 | 000112-37-8 | 72 |
| 3 | Nonanoic acid | 158 | C9H18O2 | 000112-05-0 | 58 |
| 4 | Ethanone, 1-(4,5-dihydro-2-thiaz... | 129 | C5H7NOS | 029926-41-8 | 46 |
| 5 | Octanoic acid | 144 | C8H16O2 | 000124-07-2 | 46 |



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Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

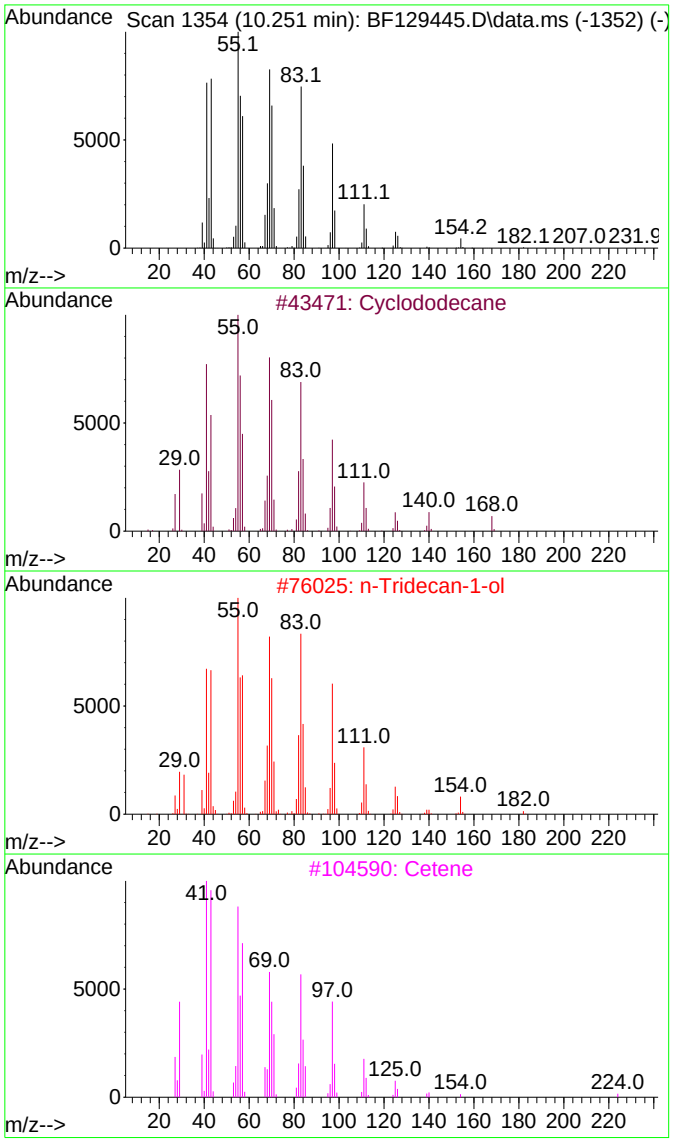
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

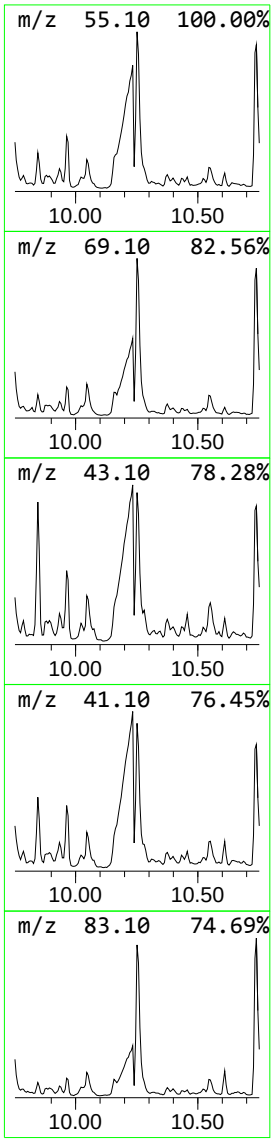
**********

Peak Number 6 Cyclododecane Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.251 | 48.29 ng | 5918090 | Acenaphthene-d10 | 9.928 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclododecane | 168 | C12H24 | 000294-62-2 | 91 |
| 2 | n-Tridecan-1-ol | 200 | C13H28O | 000112-70-9 | 91 |
| 3 | Cetene | 224 | C16H32 | 000629-73-2 | 91 |
| 4 | Formic acid, dodecyl ester | 214 | C13H26O2 | 028303-42-6 | 91 |
| 5 | Pentadecafluorooctanoic acid, do... | 582 | C20H25F15O2 | 1000406-04-2 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

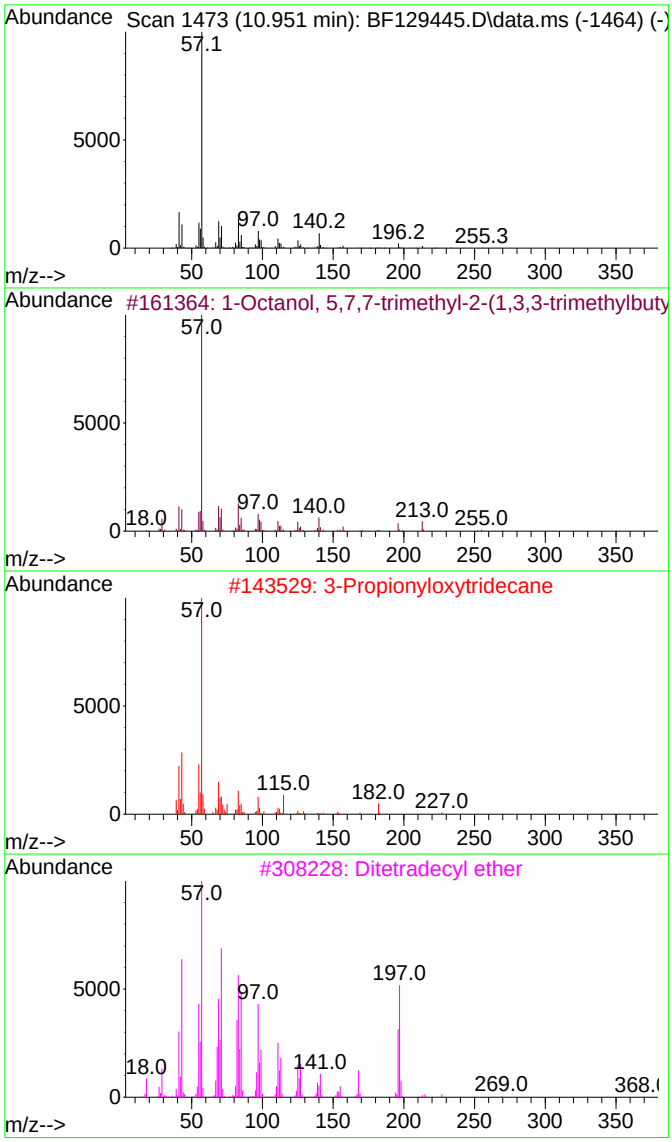
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

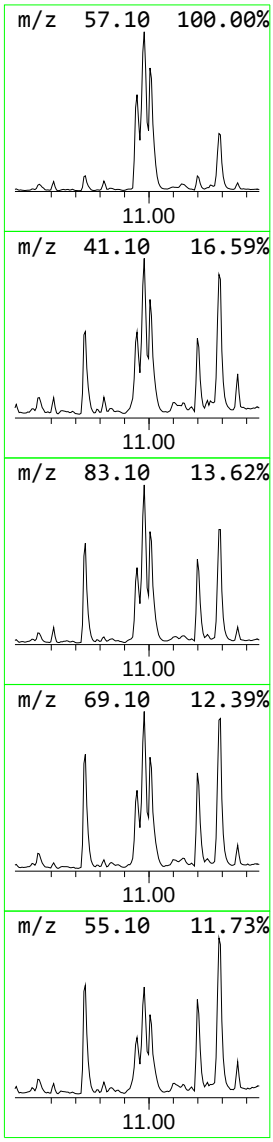
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 7 1-Octanol, 5,7,7-trimethyl-... Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.951 | 157.77 ng | 8225580 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Octanol, 5,7,7-trimethyl-2-(1,... | 270 | C18H38O | 036400-98-3 | 80 |
| 2 | 3-Propionyloxytridecane | 256 | C16H32O2 | 1000245-62-5 | 53 |
| 3 | Ditetradecyl ether | 410 | C28H58O | 005412-98-6 | 53 |
| 4 | Hexanoic acid, 3,5,5-trimethyl-,... | 340 | C22H44O2 | 1000406-26-2 | 50 |
| 5 | Pentafluoropropionic acid, tetra... | 360 | C17H29F5O2 | 006222-06-6 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

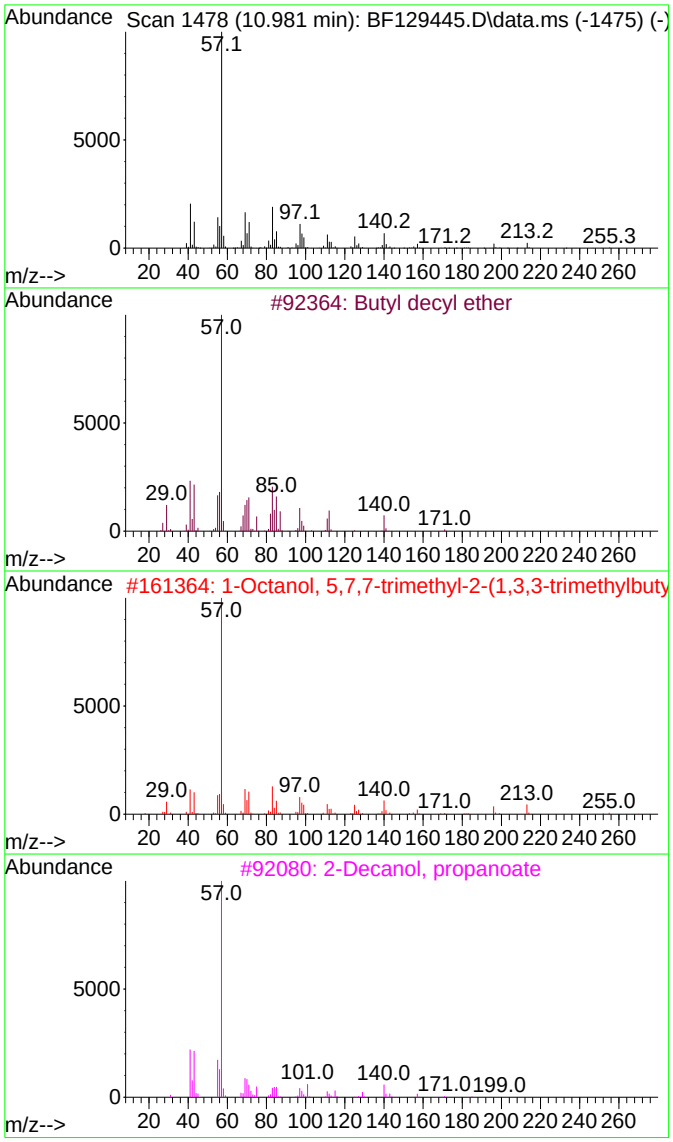
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

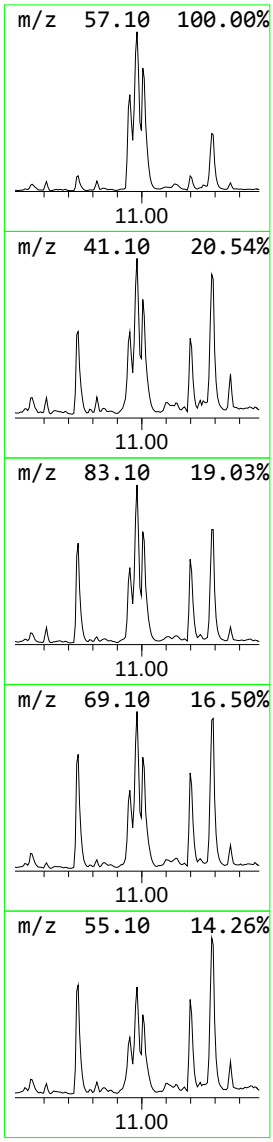
**********

Peak Number 8 Butyl decyl ether Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.981 | 271.07 ng | 14133000 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butyl decyl ether | 214 | C14H30O | 1000406-40-2 | 72 |
| 2 | 1-Octanol, 5,7,7-trimethyl-2-(1,... | 270 | C18H38O | 036400-98-3 | 72 |
| 3 | 2-Decanol, propanoate | 214 | C13H26O2 | 055683-11-9 | 53 |
| 4 | Butyl octadecyl ether | 326 | C22H46O | 1000406-40-6 | 53 |
| 5 | Tetracosyl heptafluorobutyrate | 550 | C28H49F7O2 | 1000351-83-7 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

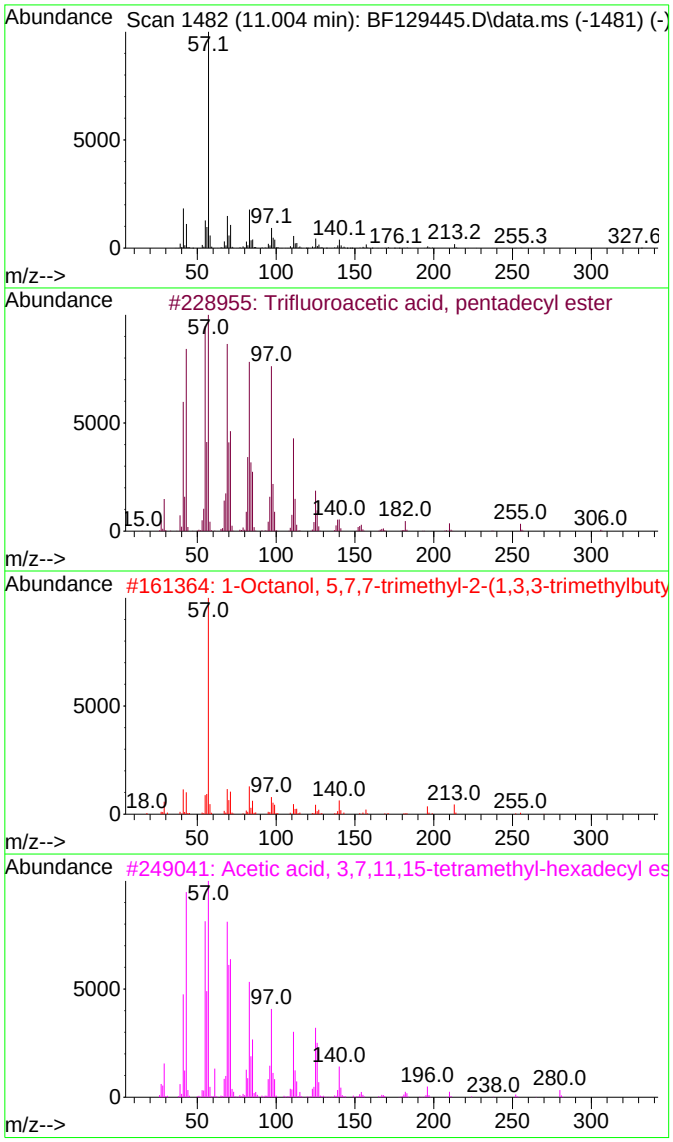
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

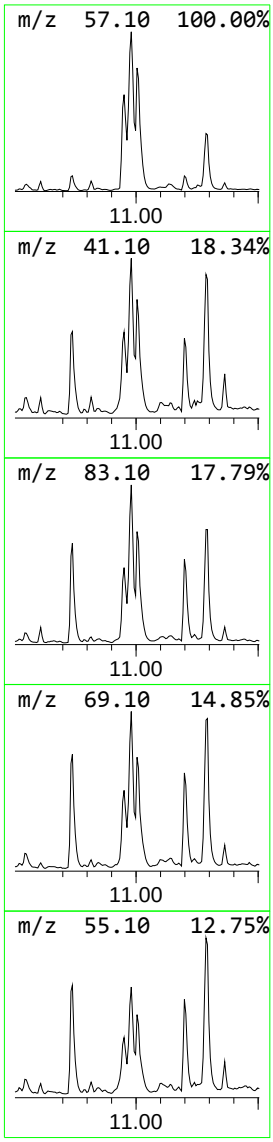
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 Trifluoroacetic acid, penta... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.004 | 168.60 ng | 8790490 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Trifluoroacetic acid, pentadecyl... | 324 | C17H31F3O2 | 959010-23-2 | 53 |
| 2 | 1-Octanol, 5,7,7-trimethyl-2-(1,... | 270 | C18H38O | 036400-98-3 | 53 |
| 3 | Acetic acid, 3,7,11,15-tetrameth... | 340 | C22H44O2 | 1000193-63-0 | 53 |
| 4 | 1-Hexacosene | 364 | C26H52 | 018835-33-1 | 53 |
| 5 | Pentafluoropropionic acid, penta... | 374 | C18H31F5O2 | 959092-08-1 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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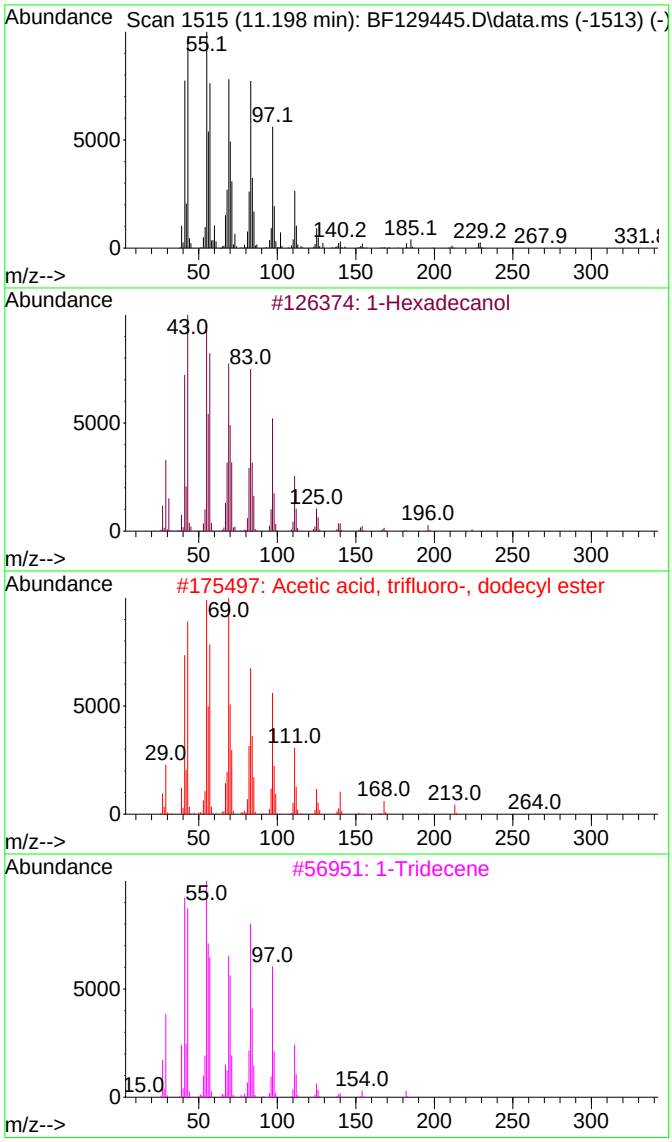
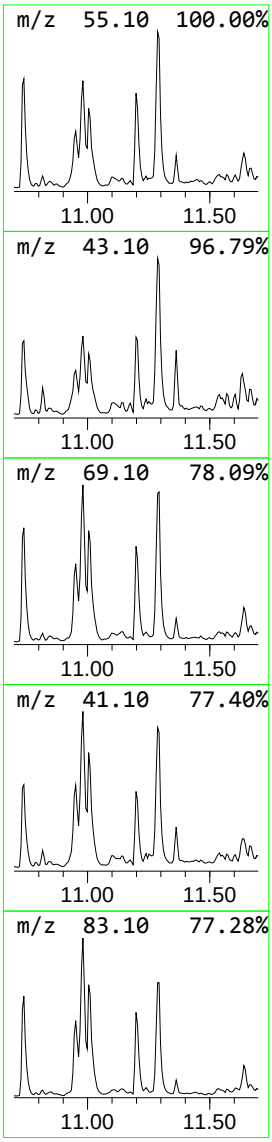
Peak Number 10 1-Hexadecanol Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.198 | 89.27 ng | 4654360 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Hexadecanol | 242 | C16H34O | 036653-82-4 | 95 |
| 2 | Acetic acid, trifluoro-, dodecyl... | 282 | C14H25F3O2 | 006222-00-0 | 94 |
| 3 | 1-Tridecene | 182 | C13H26 | 002437-56-1 | 93 |
| 4 | Cetene | 224 | C16H32 | 000629-73-2 | 91 |
| 5 | n-Pentadecanol | 228 | C15H32O | 000629-76-5 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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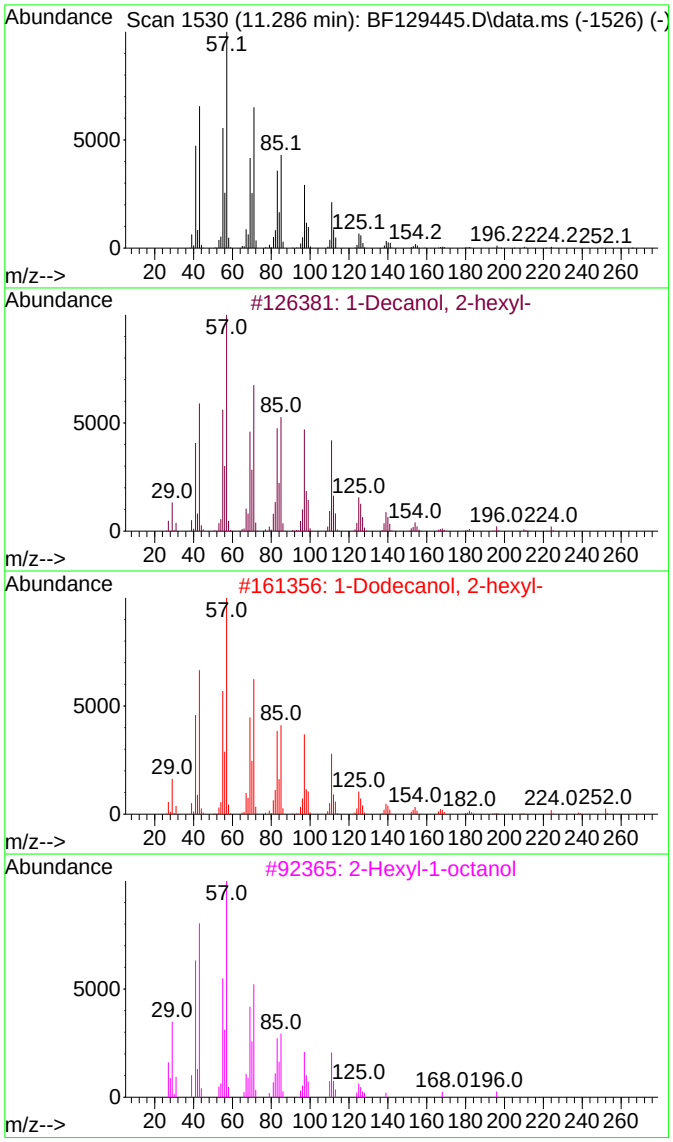
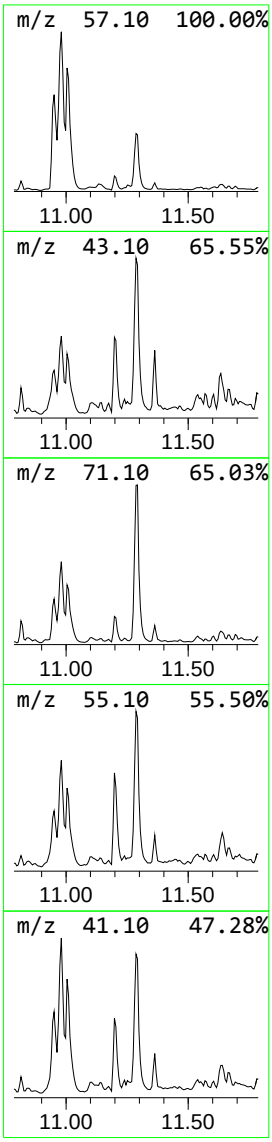
Peak Number 11 1-Decanol, 2-hexyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.286 | 229.68 ng | 11974500 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Decanol, 2-hexyl- | 242 | C16H34O | 002425-77-6 | 91 |
| 2 | 1-Dodecanol, 2-hexyl- | 270 | C18H38O | 110225-00-8 | 91 |
| 3 | 2-Hexyl-1-octanol | 214 | C14H30O | 019780-79-1 | 91 |
| 4 | 1-Dodecanol, 2-octyl- | 298 | C20H42O | 005333-42-6 | 91 |
| 5 | 1-Decanol, 2-octyl- | 270 | C18H38O | 045235-48-1 | 91 |



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Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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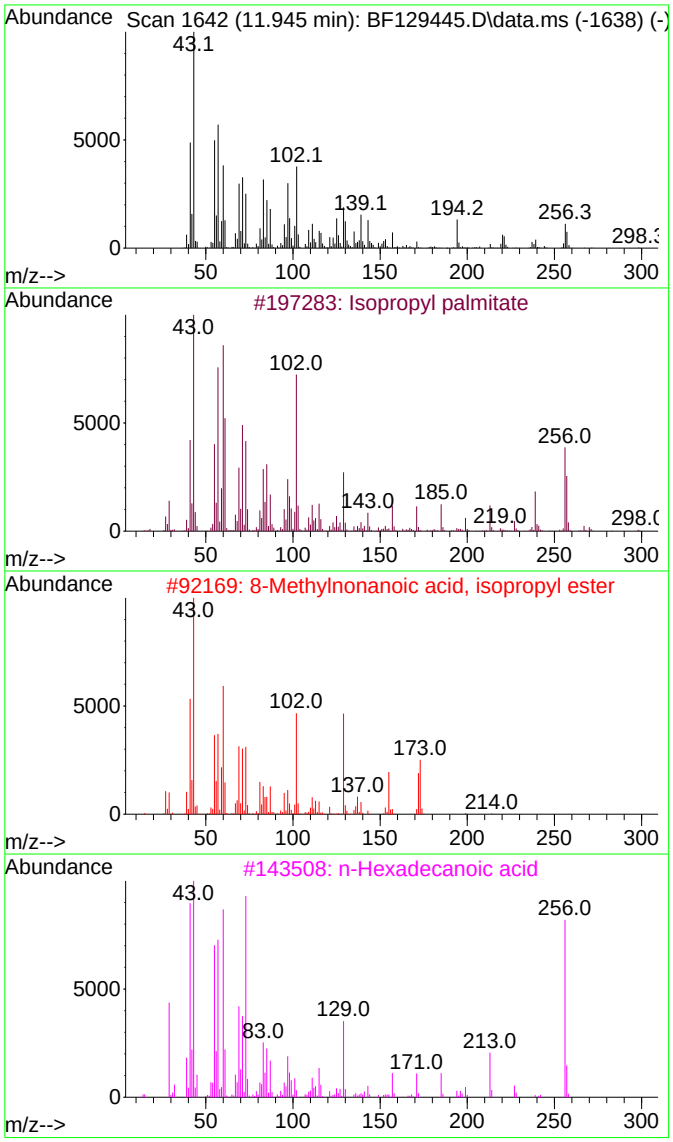
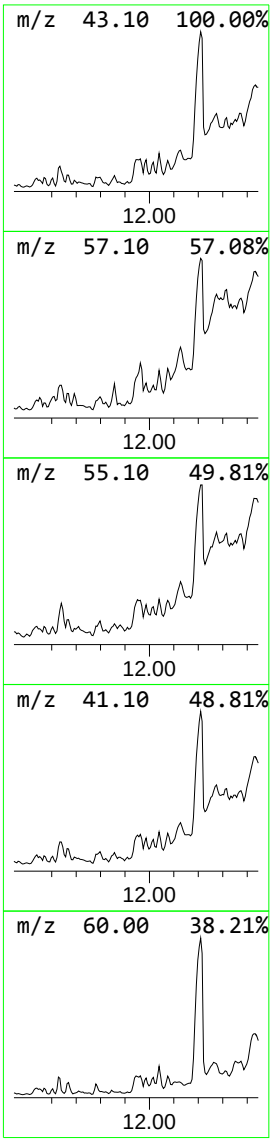
Peak Number 12 unknown11.945 Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.945 | 59.99 ng | 3127940 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Isopropyl palmitate | 298 | C19H38O2 | 000142-91-6 | 45 |
| 2 | 8-Methylnonanoic acid, isopropyl... | 214 | C13H26O2 | 1000452-00-6 | 38 |
| 3 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 35 |
| 4 | 8-Methylnonanoic acid, n-propyl ... | 214 | C13H26O2 | 1010452-00-5 | 35 |
| 5 | n-Capric acid isopropyl ester | 214 | C13H26O2 | 002311-59-3 | 27 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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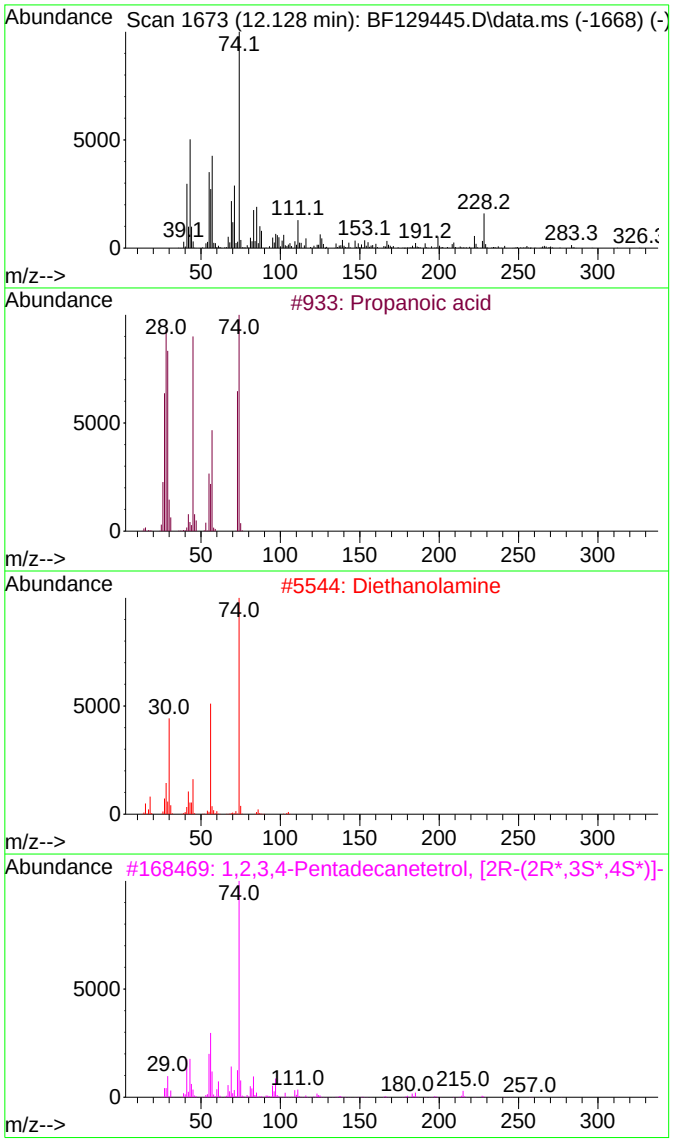
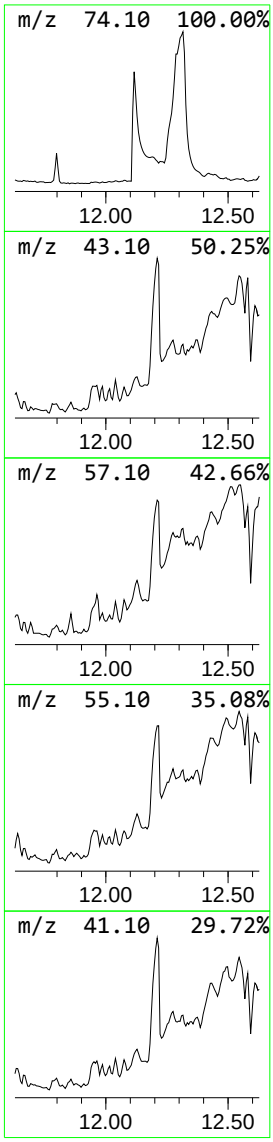
Peak Number 13 unknown12.128 Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.128 | 68.67 ng | 3580130 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propanoic acid | 74 | C3H6O2 | 000079-09-4 | 43 |
| 2 | Diethanolamine | 105 | C4H11NO2 | 000111-42-2 | 43 |
| 3 | 1,2,3,4-Pentadecanetetrol, [2R-(... | 276 | C15H32O4 | 136953-05-4 | 37 |
| 4 | Methylcyclohexylacetate | 156 | C9H16O2 | 014352-61-5 | 30 |
| 5 | Undecane, 2-methyl- | 170 | C12H26 | 007045-71-8 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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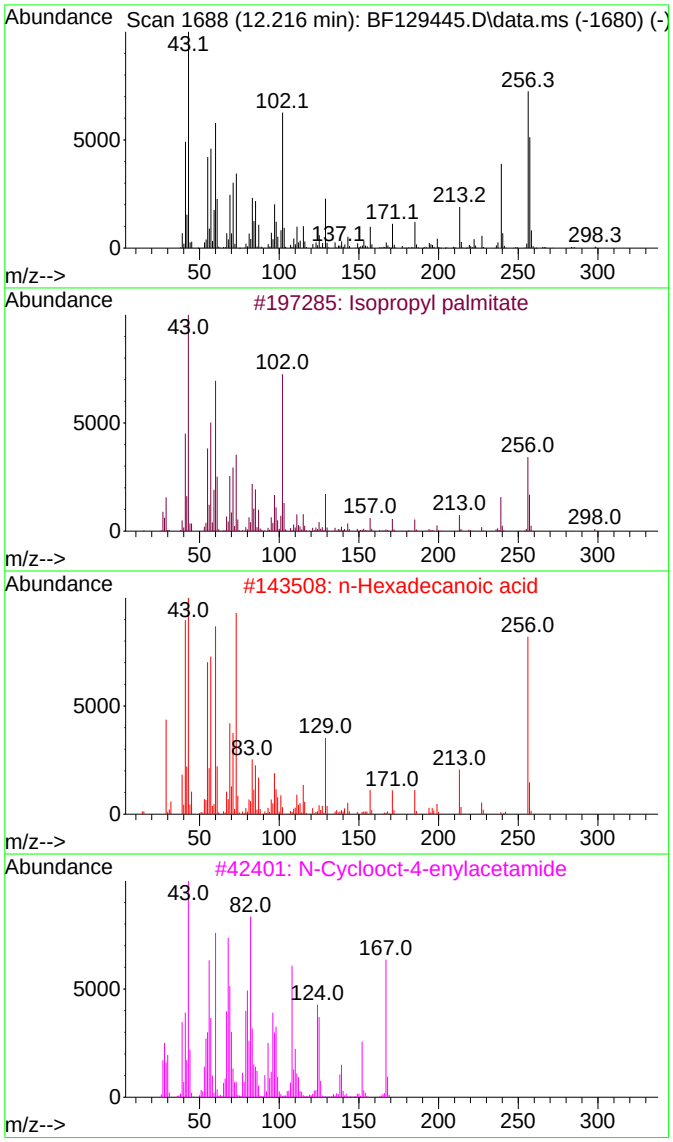
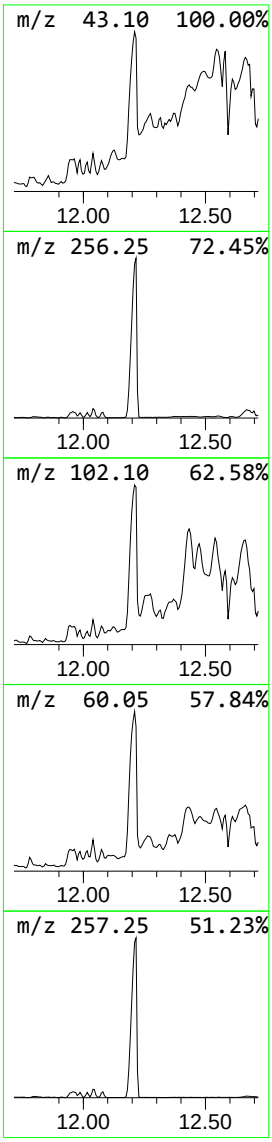
Peak Number 14 Isopropyl palmitate Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.216 | 427.11 ng | 22268100 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Isopropyl palmitate | 298 | C19H38O2 | 000142-91-6 | 95 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 60 |
| 3 | N-Cyclooct-4-enylacetamide | 167 | C10H17NO | 170952-69-9 | 25 |
| 4 | Estra-1,3,5(10)-trien-17.beta.-ol | 256 | C18H24O | 002529-64-8 | 25 |
| 5 | N-(4-Methylphenyl)methanesulfona... | 257 | C11H19NO2SSi | 1000486-49-3 | 15 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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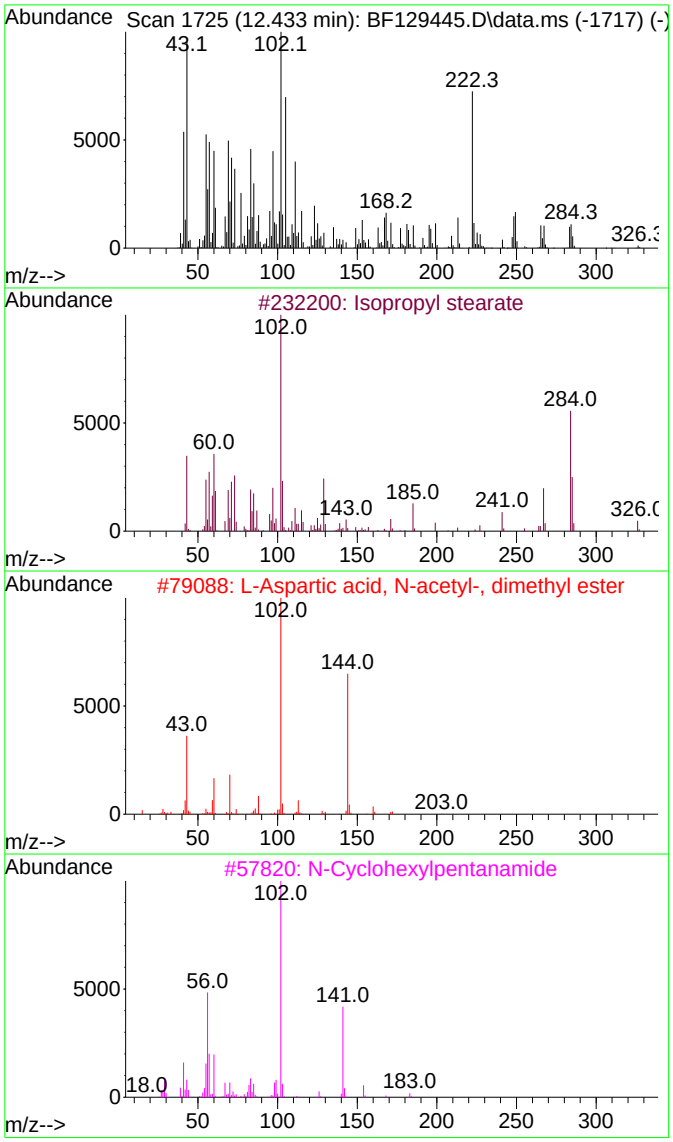
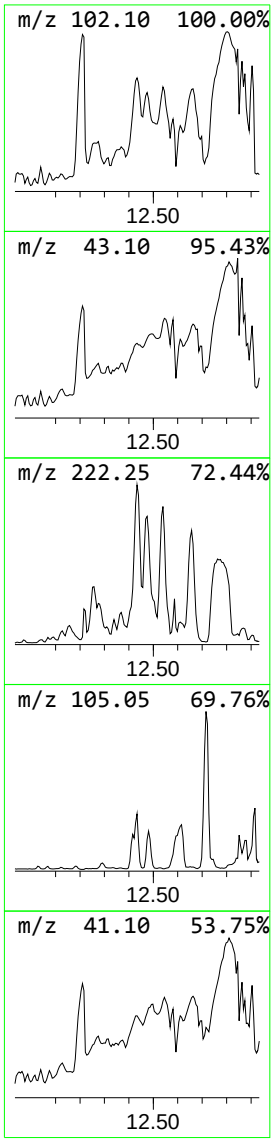
Peak Number 15 unknown12.433 Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.433 | 347.92 ng | 18139600 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Isopropyl stearate | 326 | C21H42O2 | 000112-10-7 | 41 |
| 2 | L-Aspartic acid, N-acetyl-, dime... | 203 | C8H13NO5 | 057289-64-2 | 14 |
| 3 | N-Cyclohexylpentanamide | 183 | C11H21NO | 002763-66-8 | 14 |
| 4 | 1-Heneicosanol | 312 | C21H44O | 015594-90-8 | 11 |
| 5 | Bromoacetic acid, hexadecyl ester | 362 | C18H35BrO2 | 005454-48-8 | 11 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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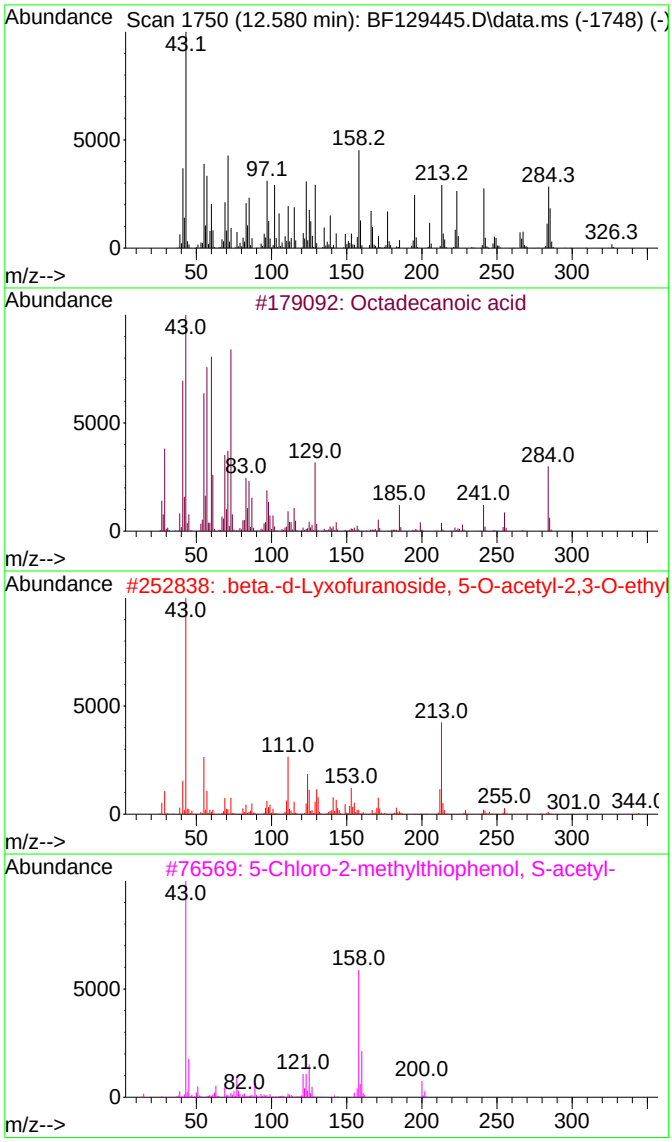
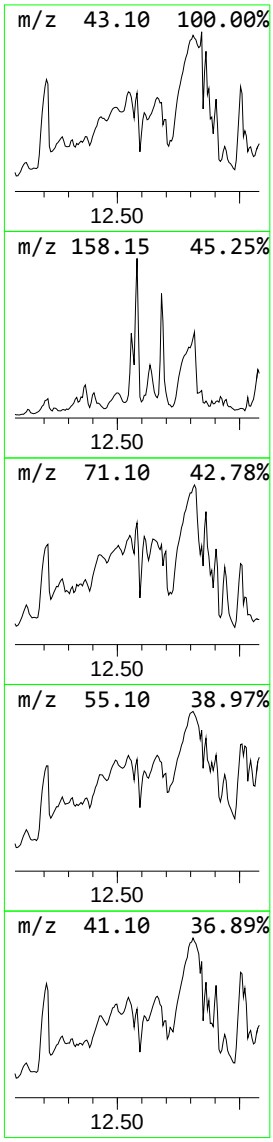
Peak Number 16 unknown12.580 Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.580 | 133.59 ng | 6965080 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 25 |
| 2 | .beta.-d-Lyxofuranoside, 5-O-ace... | 344 | C16H29BO5S | 1000154-06-5 | 10 |
| 3 | 5-Chloro-2-methylthiophenol, S-a... | 200 | C9H9ClOS | 1000463-44-0 | 10 |
| 4 | .beta.-d-Lyxofuranoside, 5-O-ace... | 372 | C18H33BO5S | 1000154-39-0 | 9 |
| 5 | 2-Chloro-6-methylthiophenol, S-a... | 200 | C9H9ClOS | 1000463-12-9 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

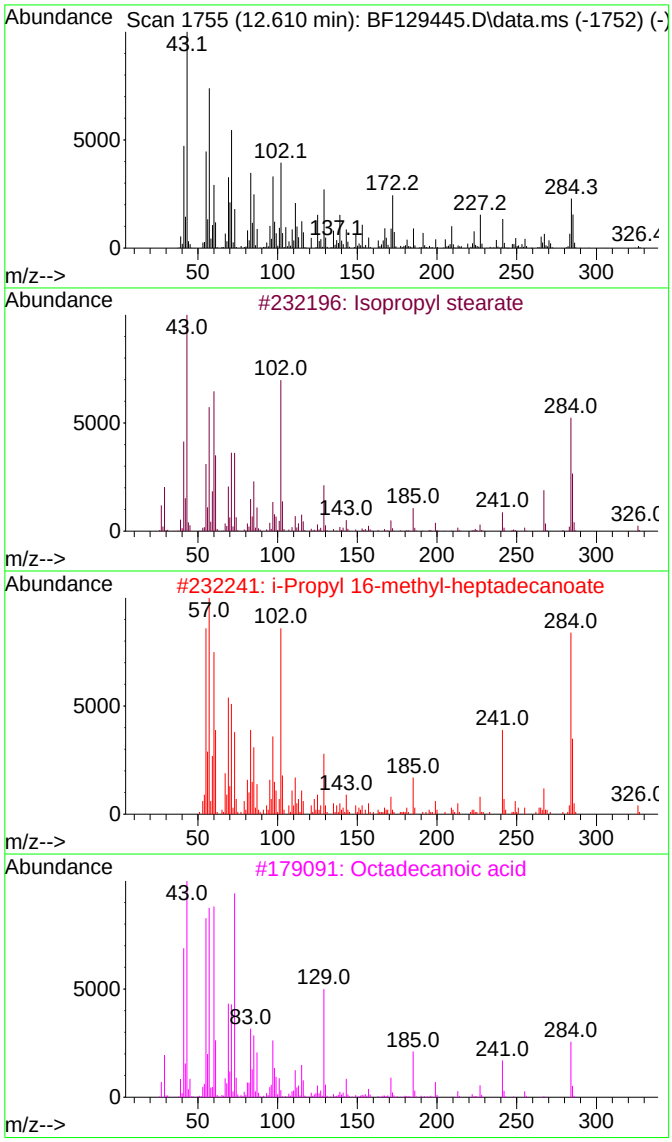
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

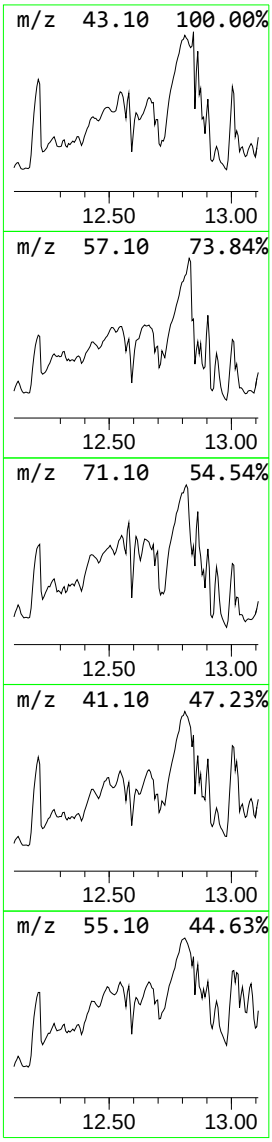
**********

Peak Number 17 unknown12.610 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.610 | 198.70 ng | 10359600 | Phenanthrene-d10 | 11.422 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Isopropyl stearate | 326 | C21H42O2 | 000112-10-7 | 42 |
| 2 | i-Propyl 16-methyl-heptadecanoate | 326 | C21H42O2 | 1000336-66-3 | 30 |
| 3 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 27 |
| 4 | n-Capric acid isopropyl ester | 214 | C13H26O2 | 002311-59-3 | 22 |
| 5 | 3,6-Epoxy-2H,8H-pyrimido[6,1-b][... | 284 | C12H16N2O6 | 041547-69-7 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

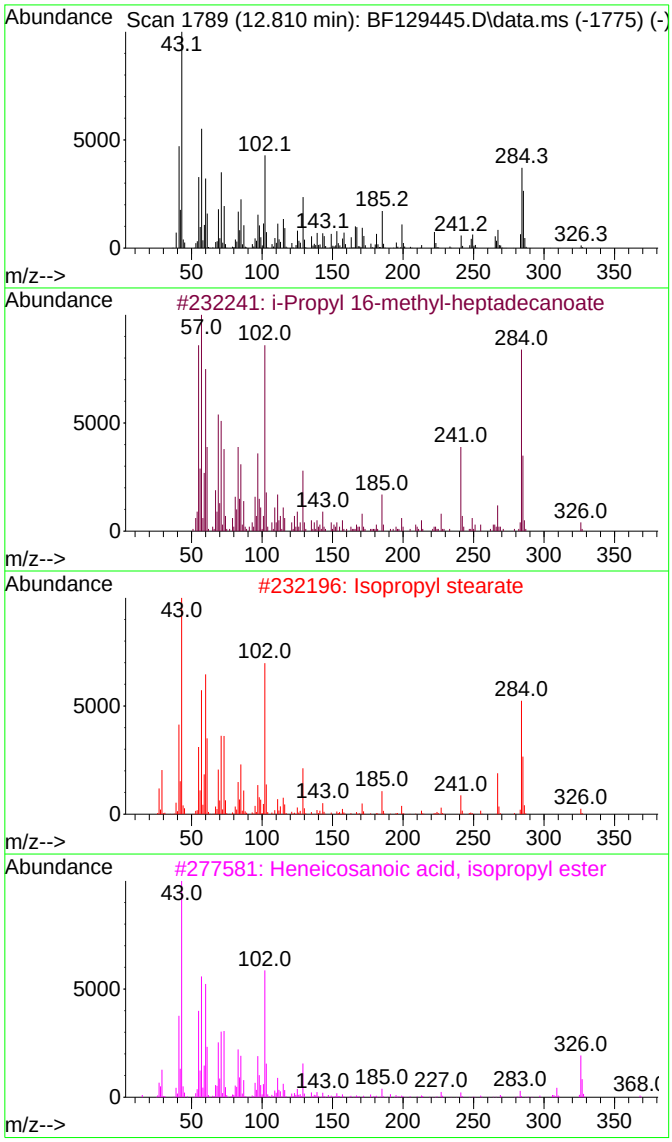
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

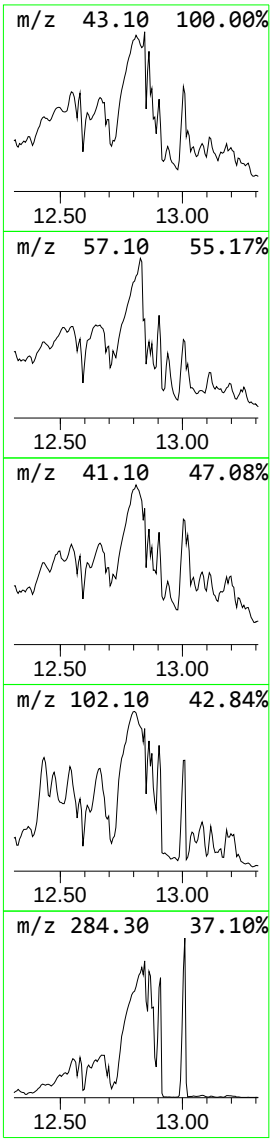
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 18 i-Propyl 16-methyl-heptadec... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.810 | 48.78 ng | 65357000 | Chrysene-d12 | 14.098 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | i-Propyl 16-methyl-heptadecanoate | 326 | C21H42O2 | 1000336-66-3 | 89 |
| 2 | Isopropyl stearate | 326 | C21H42O2 | 000112-10-7 | 46 |
| 3 | Heneicosanoic acid, isopropyl ester | 368 | C24H48O2 | 1000405-13-6 | 27 |
| 4 | Heptanoic acid, 2,2-dimethyl-6-o... | 186 | C10H18O3 | 000926-27-2 | 25 |
| 5 | 8-Methylnonanoic acid, isopropyl... | 214 | C13H26O2 | 1000452-00-6 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

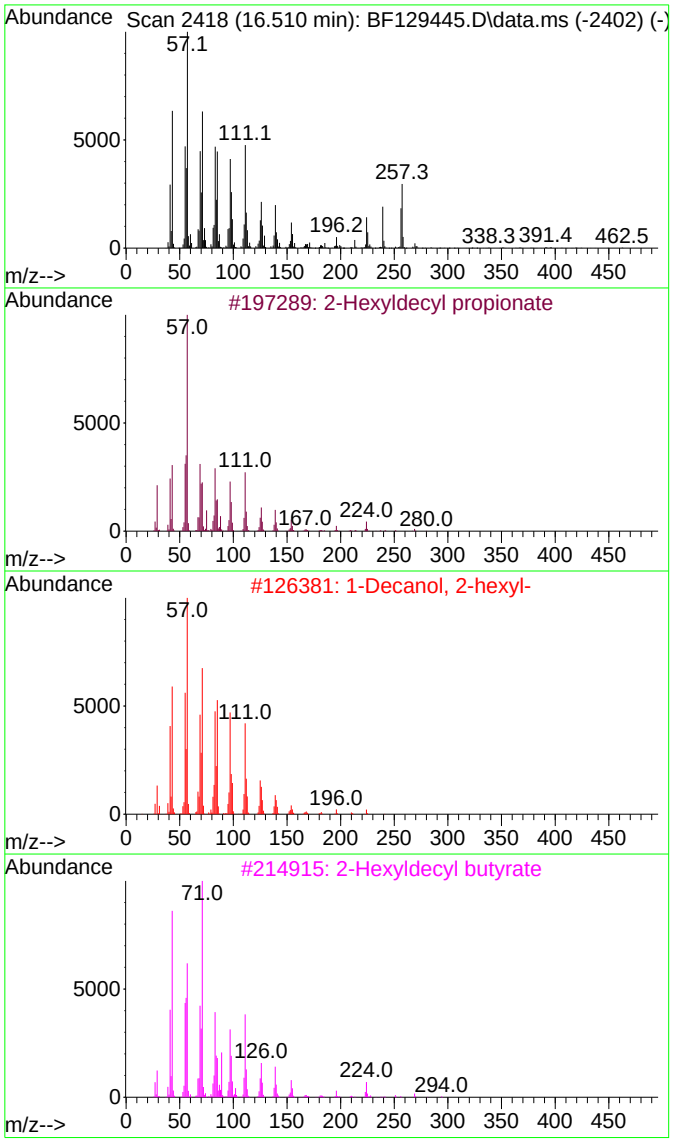
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

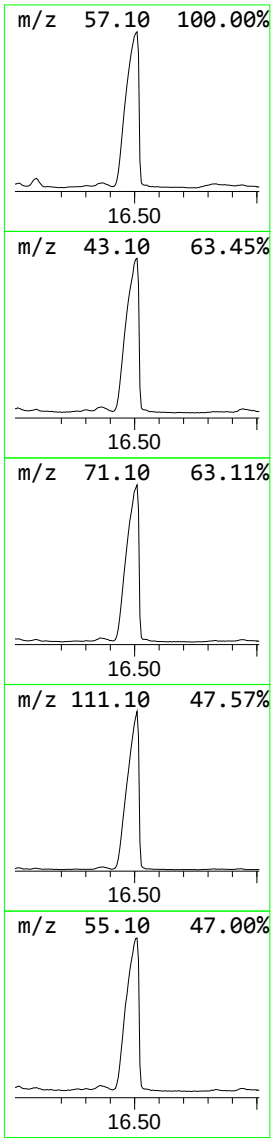
**********

Peak Number 19 2-Hexyldecyl propionate Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.510 | 737.24 ng | 39889400 | Perylene-d12 | 15.569 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Hexyldecyl propionate | 298 | C19H38O2 | 1072781-99-7 | 68 |
| 2 | 1-Decanol, 2-hexyl- | 242 | C16H34O | 002425-77-6 | 64 |
| 3 | 2-Hexyldecyl butyrate | 312 | C20H40O2 | 1072782-00-3 | 49 |
| 4 | 1-Decanol, 2-octyl- | 270 | C18H38O | 045235-48-1 | 49 |
| 5 | 1-Dodecanol, 2-hexyl- | 270 | C18H38O | 110225-00-8 | 49 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

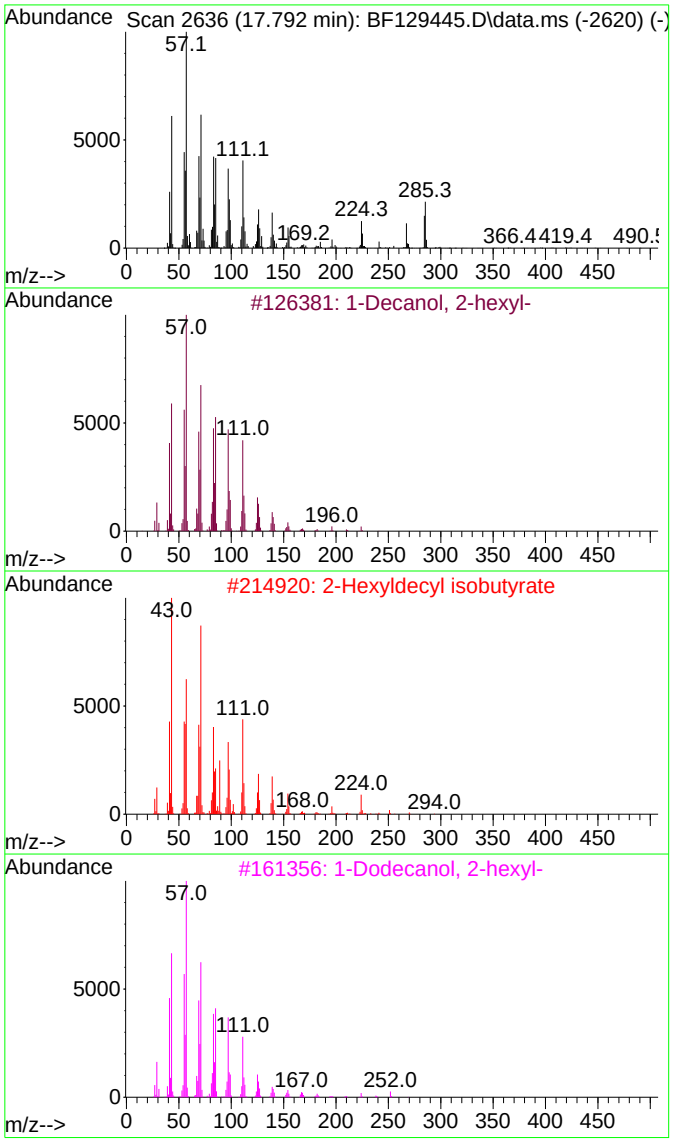
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

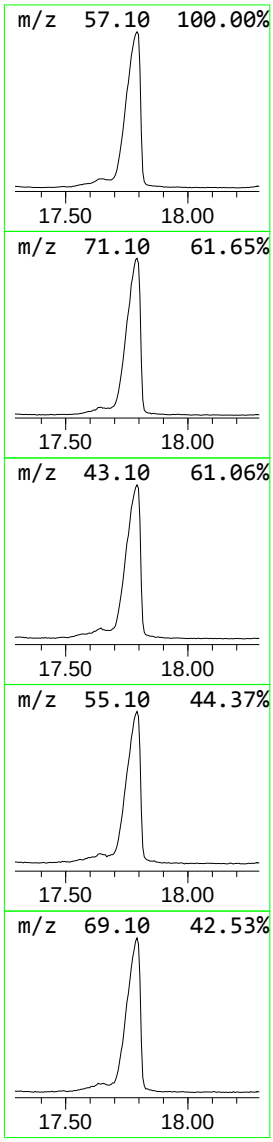
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 20 2-Hexyldecyl isobutyrate Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.792 | 574.45 ng | 31081200 | Perylene-d12 | 15.569 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Decanol, 2-hexyl- | 242 | C16H34O | 002425-77-6 | 81 |
| 2 | 2-Hexyldecyl isobutyrate | 312 | C20H40O2 | 1000368-55-3 | 76 |
| 3 | 1-Dodecanol, 2-hexyl- | 270 | C18H38O | 110225-00-8 | 68 |
| 4 | 1-Decanol, 2-octyl- | 270 | C18H38O | 045235-48-1 | 68 |
| 5 | Ethanol, 2-(hexadecyloxy)- | 286 | C18H38O2 | 002136-71-2 | 58 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072922\
Data File : BF129445.D
Acq On : 29 Jul 2022 16:58
Operator : CG\JU
Sample : N3949-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
EFF-WW

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.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 |
| Ethanol, 2-butoxy- | 5.846 | 61.3 | ng | 2732600 | 1 | 6.881 | 892328 | 20.0 |
| 1-Hexanol, 2-et... | 6.957 | 203.4 | ng | 9074090 | 1 | 6.881 | 892328 | 20.0 |
| Hexanoic acid, ... | 7.928 | 991.7 | ng | 71004900 | 2 | 8.169 | 1431940 | 20.0 |
| n-Decanoic acid | 9.187 | 68.0 | ng | 8333480 | 3 | 9.928 | 2451070 | 20.0 |
| Dodecanoic acid | 10.233 | 150.1 | ng | 18390400 | 3 | 9.928 | 2451070 | 20.0 |
| Cyclododecane | 10.251 | 48.3 | ng | 5918090 | 3 | 9.928 | 2451070 | 20.0 |
| 1-Octanol, 5,7,... | 10.951 | 157.8 | ng | 8225580 | 4 | 11.422 | 1042740 | 20.0 |
| Butyl decyl ether | 10.981 | 271.1 | ng | 14133000 | 4 | 11.422 | 1042740 | 20.0 |
| Trifluoroacetic... | 11.004 | 168.6 | ng | 8790490 | 4 | 11.422 | 1042740 | 20.0 |
| 1-Hexadecanol | 11.198 | 89.3 | ng | 4654360 | 4 | 11.422 | 1042740 | 20.0 |
| 1-Decanol, 2-he... | 11.286 | 229.7 | ng | 11974500 | 4 | 11.422 | 1042740 | 20.0 |
| unknown11.945 | 11.945 | 60.0 | ng | 3127940 | 4 | 11.422 | 1042740 | 20.0 |
| unknown12.128 | 12.128 | 68.7 | ng | 3580130 | 4 | 11.422 | 1042740 | 20.0 |
| Isopropyl palmi... | 12.216 | 427.1 | ng | 22268100 | 4 | 11.422 | 1042740 | 20.0 |
| unknown12.433 | 12.433 | 347.9 | ng | 18139600 | 4 | 11.422 | 1042740 | 20.0 |
| unknown12.580 | 12.580 | 133.6 | ng | 6965080 | 4 | 11.422 | 1042740 | 20.0 |
| unknown12.610 | 12.610 | 198.7 | ng | 10359600 | 4 | 11.422 | 1042740 | 20.0 |
| i-Propyl 16-met... | 12.810 | 48.8 | ng | 65357000 | 5 | 14.098 | 26795100 | 20.0 |
| 2-Hexyldecyl pr... | 16.510 | 737.2 | ng | 39889400 | 6 | 15.569 | 1082120 | 20.0 |
| 2-Hexyldecyl is... | 17.792 | 574.5 | ng | 31081200 | 6 | 15.569 | 1082120 | 20.0 |