LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

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-BF042123.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF132978.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 2.540 | 70 | 77 | 91 | rVB | 700739 | 1186069 | 16.64% | 1.656% |
| 2 | 4.945 | 481 | 486 | 493 | rVB | 1504157 | 1631887 | 22.90% | 2.279% |
| 3 | 5.222 | 529 | 533 | 539 | rVB | 135121 | 118581 | 1.66% | 0.166% |
| 4 | 5.351 | 548 | 555 | 559 | rBV | 2376745 | 2550516 | 35.79% | 3.561% |
| 5 | 5.540 | 582 | 587 | 590 | rBV | 147034 | 129640 | 1.82% | 0.181% |
| 6 | 5.592 | 592 | 596 | 599 | rBV | 400550 | 359409 | 5.04% | 0.502% |
| 7 | 5.757 | 616 | 624 | 626 | rVB | 45974 | 71511 | 1.00% | 0.100% |
| 8 | 6.051 | 671 | 674 | 683 | rVB | 547873 | 468346 | 6.57% | 0.654% |
| 9 | 6.363 | 721 | 727 | 730 | rBV | 103368 | 125340 | 1.76% | 0.175% |
| 10 | 6.398 | 730 | 733 | 741 | rVB | 873581 | 1379540 | 19.36% | 1.926% |
| 11 | 6.569 | 755 | 762 | 764 | rBV | 2906221 | 3845408 | 53.96% | 5.369% |
| 12 | 6.598 | 764 | 767 | 770 | rVV | 3471654 | 3776177 | 52.99% | 5.273% |
| 13 | 6.622 | 770 | 771 | 777 | rVB | 141348 | 108279 | 1.52% | 0.151% |
| 14 | 6.692 | 777 | 783 | 786 | rBV | 4314665 | 4966843 | 69.69% | 6.935% |
| 15 | 6.751 | 789 | 793 | 796 | rVV | 1471410 | 1258781 | 17.66% | 1.758% |
| 16 | 6.787 | 796 | 799 | 802 | rVB | 425655 | 371763 | 5.22% | 0.519% |
| 17 | 7.151 | 857 | 861 | 868 | rVB | 152640 | 162115 | 2.27% | 0.226% |
| 18 | 7.316 | 878 | 889 | 892 | rBV | 4172056 | 4787963 | 67.18% | 6.685% |
| 19 | 7.563 | 923 | 931 | 934 | rBV6 | 37523 | 80807 | 1.13% | 0.113% |
| 20 | 7.698 | 950 | 954 | 957 | rBV2 | 82968 | 105854 | 1.49% | 0.148% |
| 21 | 7.981 | 986 | 1002 | 1005 | rBV | 1811326 | 4711648 | 66.11% | 6.579% |
| 22 | 8.034 | 1007 | 1011 | 1014 | rVB | 2172018 | 1697315 | 23.82% | 2.370% |
| 23 | 8.263 | 1046 | 1050 | 1052 | rBV | 468482 | 398691 | 5.59% | 0.557% |
| 24 | 8.286 | 1052 | 1054 | 1056 | rVV | 413269 | 378398 | 5.31% | 0.528% |
| 25 | 8.304 | 1056 | 1057 | 1062 | rVB | 179016 | 151183 | 2.12% | 0.211% |
| 26 | 8.375 | 1066 | 1069 | 1072 | rBV2 | 138681 | 117752 | 1.65% | 0.164% |
| 27 | 8.886 | 1153 | 1156 | 1163 | rVB2 | 71391 | 104648 | 1.47% | 0.146% |
| 28 | 9.116 | 1185 | 1195 | 1198 | rBV | 7000285 | 7126768 | 100.00% | 9.951% |
| 29 | 9.245 | 1213 | 1217 | 1219 | rVB2 | 143609 | 153132 | 2.15% | 0.214% |
| 30 | 9.704 | 1290 | 1295 | 1299 | rBV | 138512 | 154445 | 2.17% | 0.216% |
| 31 | 9.763 | 1301 | 1305 | 1306 | rVV | 570290 | 507156 | 7.12% | 0.708% |
| 32 | 9.786 | 1306 | 1309 | 1312 | rVV | 2277989 | 1836023 | 25.76% | 2.564% |
| 33 | 9.816 | 1312 | 1314 | 1315 | rVV | 200900 | 157241 | 2.21% | 0.220% |
| 34 | 9.839 | 1315 | 1318 | 1320 | rVV | 362121 | 359722 | 5.05% | 0.502% |
| 35 | 9.863 | 1320 | 1322 | 1327 | rVB | 234796 | 239251 | 3.36% | 0.334% |
| 36 | 10.245 | 1385 | 1387 | 1389 | rBV | 106896 | 104345 | 1.46% | 0.146% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

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-BF042123.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 10.351 | 1400 | 1405 | 1407 | rBV2 | 53244 | 98430 | 1.38% | 0.137% |
| 38 | 10.427 | 1413 | 1418 | 1421 | rBV | 3607468 | 3454560 | 48.47% | 4.824% |
| 39 | 10.539 | 1434 | 1437 | 1439 | rBV | 115276 | 82945 | 1.16% | 0.116% |
| 40 | 10.575 | 1439 | 1443 | 1446 | rBV | 3160620 | 2833758 | 39.76% | 3.957% |
| 41 | 11.033 | 1519 | 1521 | 1528 | rBV | 181961 | 192218 | 2.70% | 0.268% |
| 42 | 11.116 | 1531 | 1535 | 1537 | rVV | 65775 | 104431 | 1.47% | 0.146% |
| 43 | 11.139 | 1537 | 1539 | 1542 | rVB | 159636 | 140367 | 1.97% | 0.196% |
| 44 | 11.269 | 1557 | 1561 | 1564 | rVB | 1926059 | 1688974 | 23.70% | 2.358% |
| 45 | 11.463 | 1586 | 1594 | 1595 | rBV2 | 79858 | 188150 | 2.64% | 0.263% |
| 46 | 11.804 | 1649 | 1652 | 1660 | rBV | 627416 | 626852 | 8.80% | 0.875% |
| 47 | 12.080 | 1696 | 1699 | 1704 | rVB3 | 58936 | 73978 | 1.04% | 0.103% |
| 48 | 12.186 | 1714 | 1717 | 1721 | rBV | 316211 | 351698 | 4.93% | 0.491% |
| 49 | 12.527 | 1764 | 1775 | 1778 | rBV2 | 174319 | 526969 | 7.39% | 0.736% |
| 50 | 12.592 | 1781 | 1786 | 1797 | rVB2 | 607883 | 787819 | 11.05% | 1.100% |
| 51 | 12.857 | 1826 | 1831 | 1834 | rBV | 5260995 | 5244542 | 73.59% | 7.323% |
| 52 | 13.198 | 1886 | 1889 | 1897 | rBV | 475886 | 550229 | 7.72% | 0.768% |
| 53 | 13.474 | 1920 | 1936 | 1957 | rBV | 274314 | 1563546 | 21.94% | 2.183% |
| 54 | 13.768 | 1983 | 1986 | 1996 | rVB2 | 93228 | 131301 | 1.84% | 0.183% |
| 55 | 13.898 | 2004 | 2008 | 2012 | rVB | 1236425 | 1112803 | 15.61% | 1.554% |
| 56 | 13.980 | 2019 | 2022 | 2023 | rBV | 100058 | 91975 | 1.29% | 0.128% |
| 57 | 14.004 | 2023 | 2026 | 2038 | rVV | 654707 | 677935 | 9.51% | 0.947% |
| 58 | 14.104 | 2038 | 2043 | 2055 | rVV | 604940 | 774366 | 10.87% | 1.081% |
| 59 | 14.304 | 2067 | 2077 | 2090 | rBV | 216370 | 923061 | 12.95% | 1.289% |
| 60 | 14.945 | 2182 | 2186 | 2195 | rBV | 640776 | 1026741 | 14.41% | 1.434% |
| 61 | 15.327 | 2246 | 2251 | 2256 | rVV | 894048 | 1045163 | 14.67% | 1.459% |
| 62 | 15.380 | 2257 | 2260 | 2268 | rVB | 70652 | 92230 | 1.29% | 0.129% |
| 63 | 16.039 | 2365 | 2372 | 2383 | rBV | 387792 | 885244 | 12.42% | 1.236% |
| 64 | 16.274 | 2410 | 2412 | 2430 | rVB3 | 80232 | 198886 | 2.79% | 0.278% |
| 65 | 17.662 | 2640 | 2648 | 2663 | rBV | 126907 | 466456 | 6.55% | 0.651% |

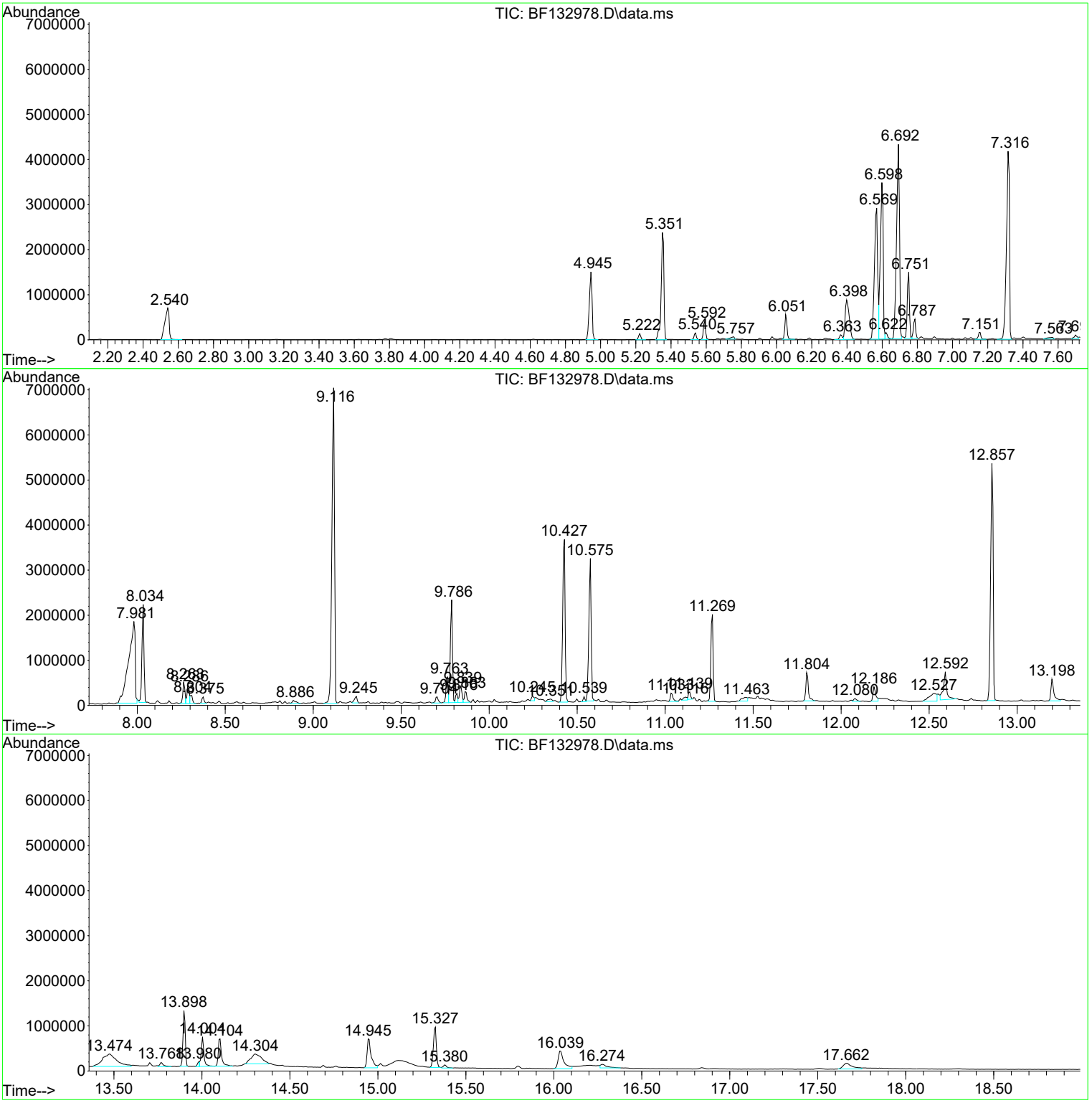
Sum of corrected areas: 71618174

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.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\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

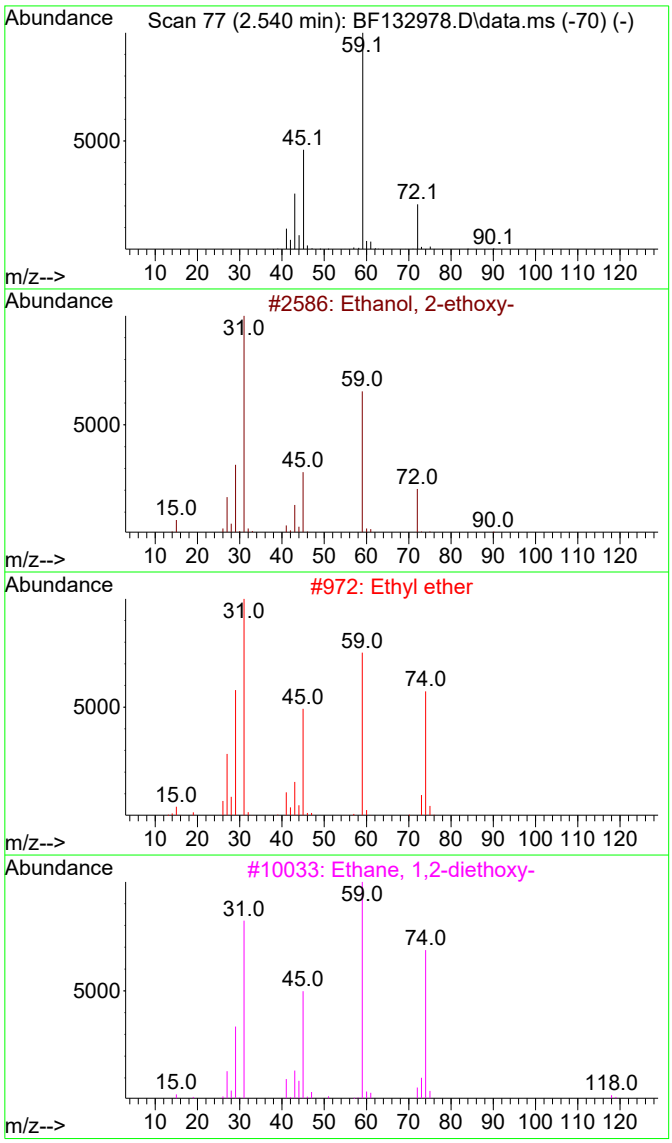
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

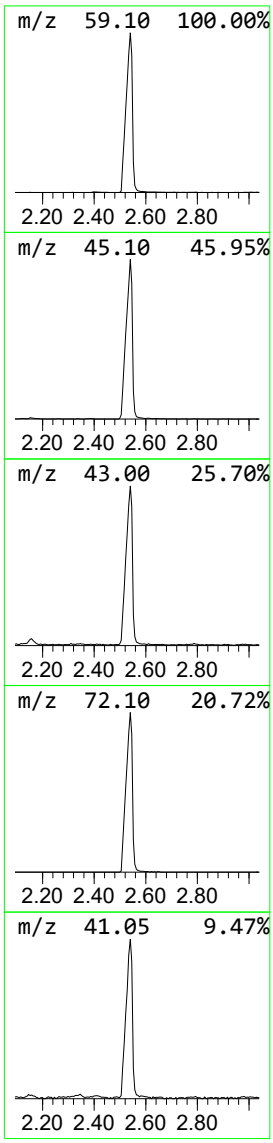
**********

Peak Number 1 Ethanol, 2-ethoxy- Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.540 | 18.84 ng | 1186070 | 1,4-Dichlorobenzene-d4 | 6.751 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Ethanol, 2-ethoxy- | 90 | C4H10O2 | 000110-80-5 | 83 |
| 2 | Ethyl ether | 74 | C4H10O | 000060-29-7 | 50 |
| 3 | Ethane, 1,2-diethoxy- | 118 | C6H14O2 | 000629-14-1 | 50 |
| 4 | Ethane, 1,1'-oxybis[2-methoxy- | 134 | C6H14O3 | 000111-96-6 | 39 |
| 5 | Hexanamide | 115 | C6H13NO | 000628-02-4 | 33 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

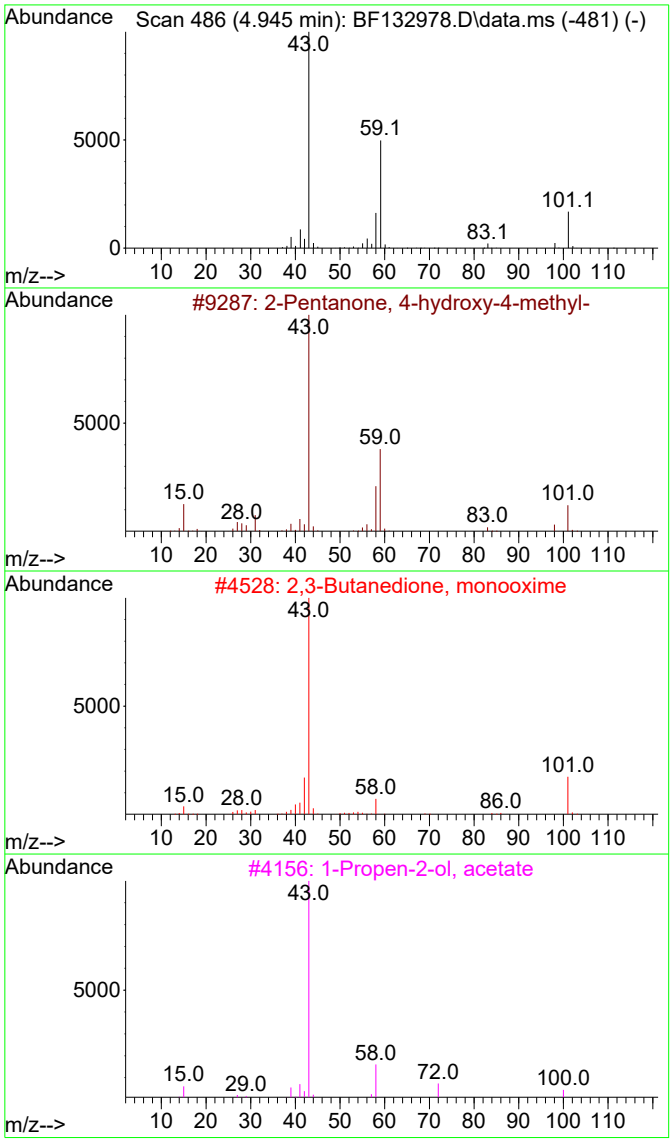
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

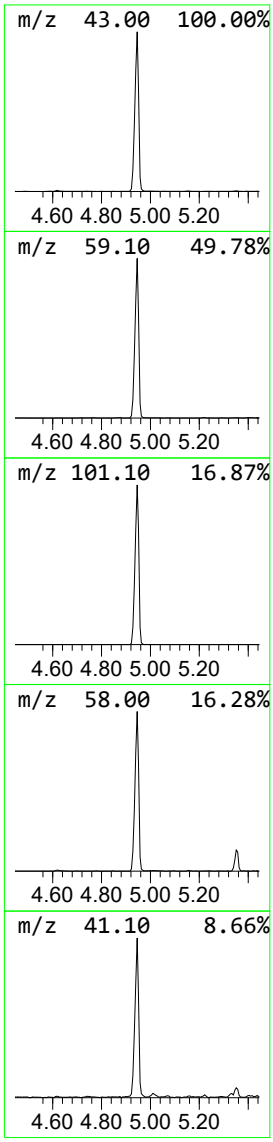
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.945 | 25.93 ng | 1631890 | 1,4-Dichlorobenzene-d4 | 6.751 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 59 |
| 2 | 2,3-Butanedione, monooxime | 101 | C4H7NO2 | 000057-71-6 | 16 |
| 3 | 1-Propen-2-ol, acetate | 100 | C5H8O2 | 000108-22-5 | 12 |
| 4 | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 9 |
| 5 | Acetone | 58 | C3H6O | 000067-64-1 | 7 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

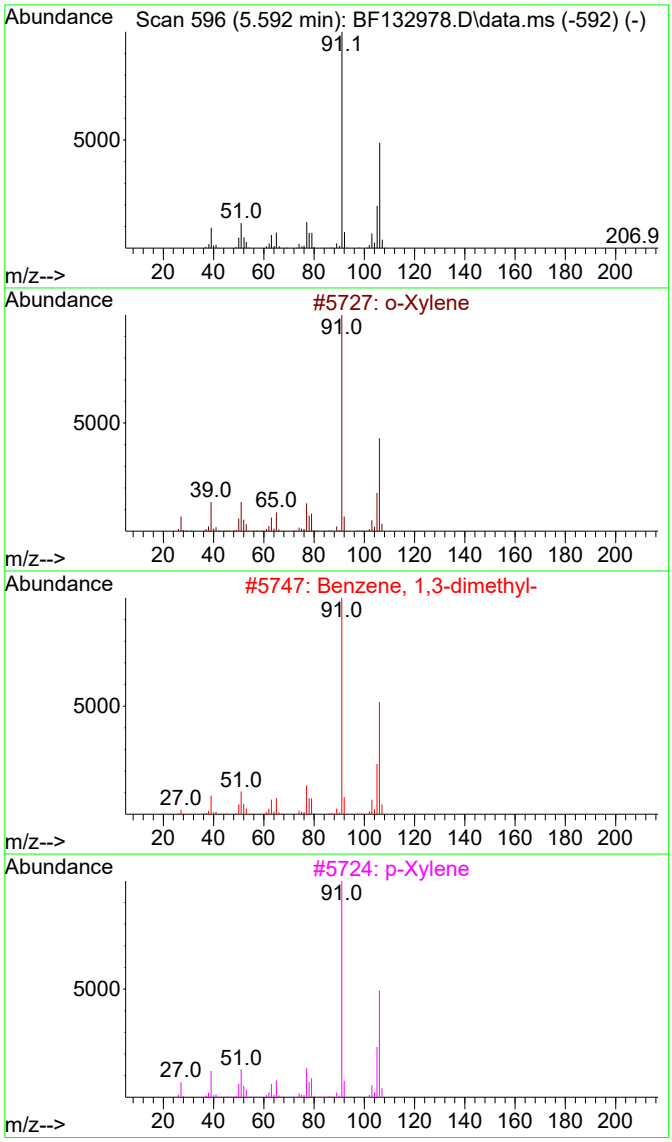
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

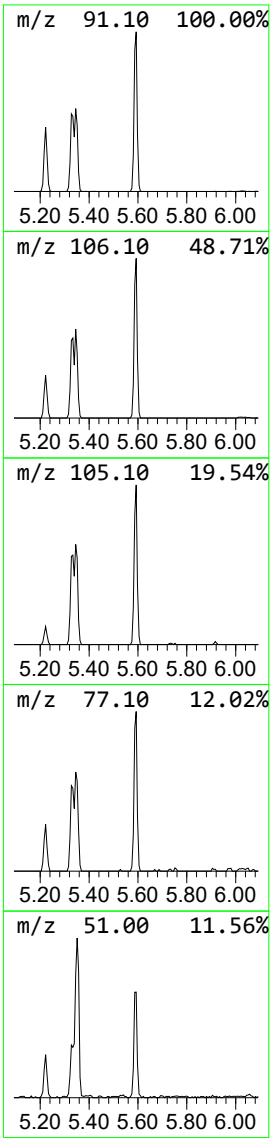
**********

Peak Number 3 o-Xylene Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.592 | 5.71 ng | 359409 | 1,4-Dichlorobenzene-d4 | 6.751 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | o-Xylene | 106 | C8H10 | 000095-47-6 | 97 |
| 2 | Benzene, 1,3-dimethyl- | 106 | C8H10 | 000108-38-3 | 97 |
| 3 | p-Xylene | 106 | C8H10 | 000106-42-3 | 97 |
| 4 | Cyclopentene, 1-ethenyl-3-methyl... | 106 | C8H10 | 061142-07-2 | 86 |
| 5 | Ethylbenzene | 106 | C8H10 | 000100-41-4 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

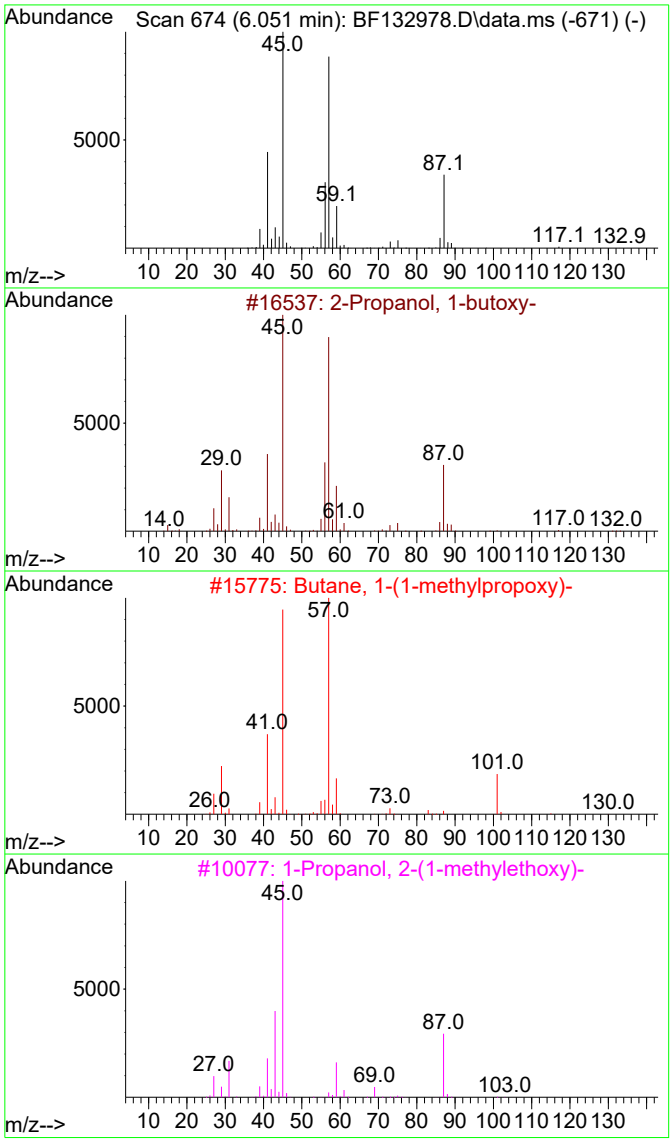
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

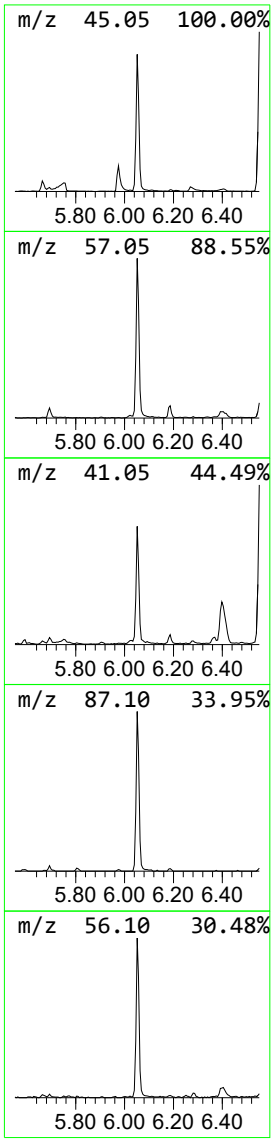
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 4 2-Propanol, 1-butoxy- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.051 | 7.44 ng | 468346 | 1,4-Dichlorobenzene-d4 | 6.751 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Propanol, 1-butoxy- | 132 | C7H16O2 | 005131-66-8 | 83 |
| 2 | Butane, 1-(1-methylpropoxy)- | 130 | C8H18O | 000999-65-5 | 53 |
| 3 | 1-Propanol, 2-(1-methylethoxy)- | 118 | C6H14O2 | 003944-37-4 | 53 |
| 4 | 2-Butanol | 74 | C4H10O | 000078-92-2 | 47 |
| 5 | Formic acid, 1-methylpropyl ester | 102 | C5H10O2 | 000589-40-2 | 45 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

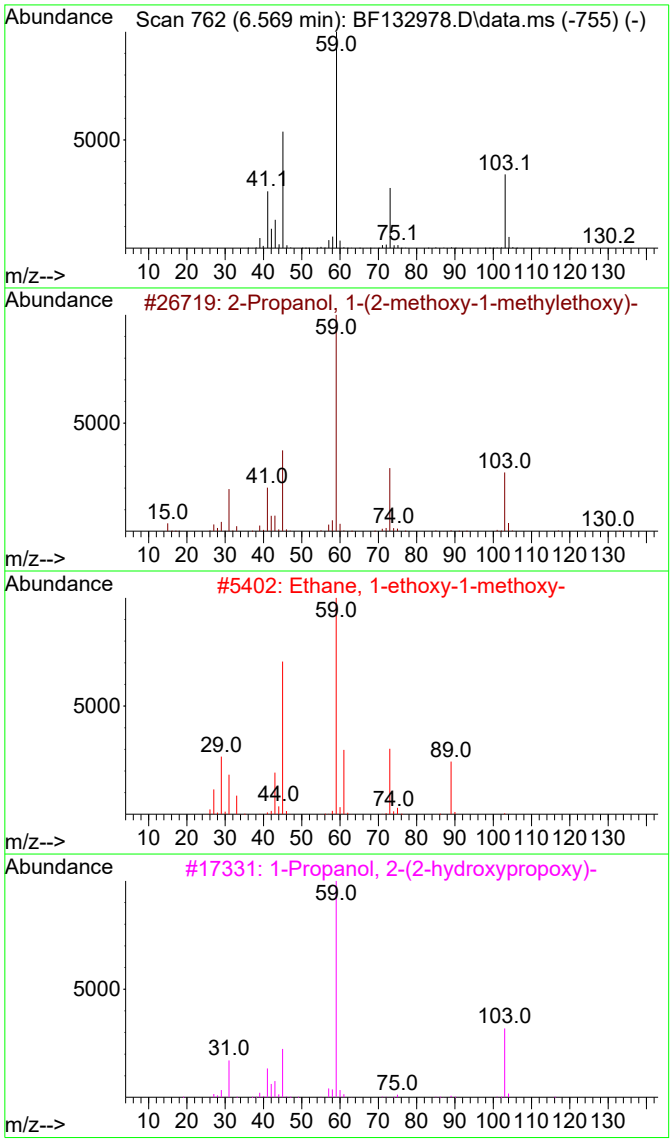
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

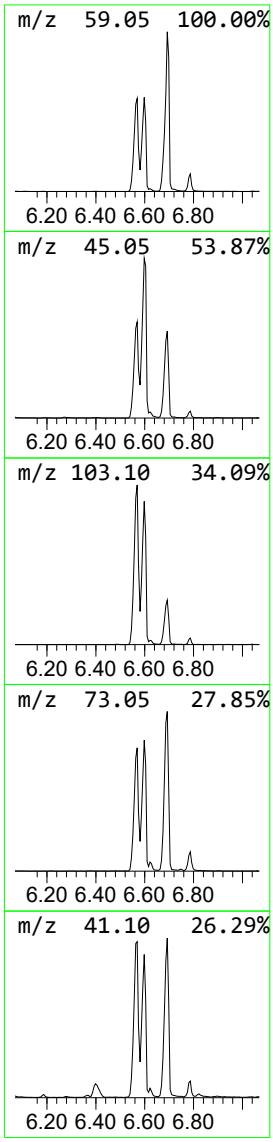
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.569 | 61.10 ng | 3845410 | 1,4-Dichlorobenzene-d4 | 6.751 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Propanol, 1-(2-methoxy-1-methy... | 148 | C7H16O3 | 020324-32-7 | 72 |
| 2 | Ethane, 1-ethoxy-1-methoxy- | 104 | C5H12O2 | 010471-14-4 | 64 |
| 3 | 1-Propanol, 2-(2-hydroxypropoxy)- | 134 | C6H14O3 | 000106-62-7 | 59 |
| 4 | 2-Propanol, 1-(2-ethoxypropoxy)- | 162 | C8H18O3 | 010143-32-5 | 56 |
| 5 | 1-Propanol, 2,2'-oxybis- | 134 | C6H14O3 | 000108-61-2 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

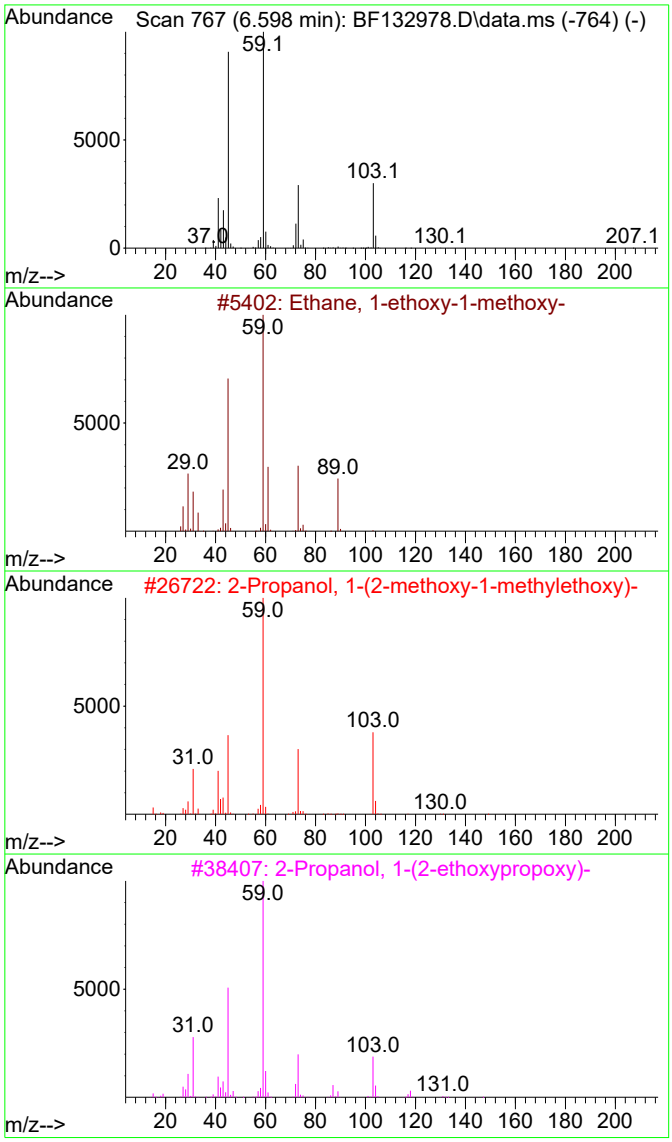
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

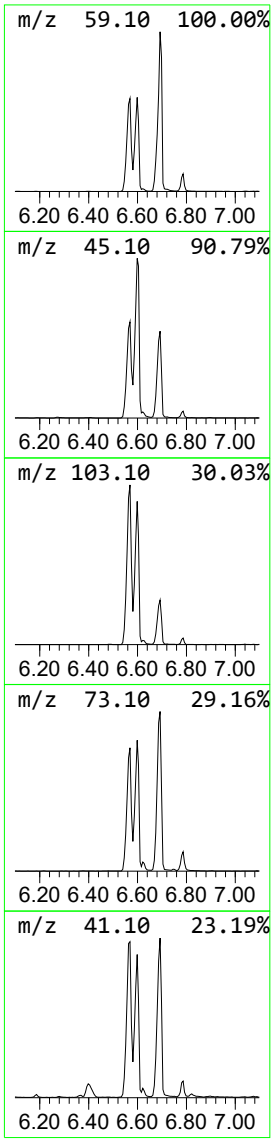
**********

Peak Number 6 Ethane, 1-ethoxy-1-methoxy- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.598 | 60.00 ng | 3776180 | 1,4-Dichlorobenzene-d4 | 6.751 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Ethane, 1-ethoxy-1-methoxy- | 104 | C5H12O2 | 010471-14-4 | 64 |
| 2 | 2-Propanol, 1-(2-methoxy-1-methy... | 148 | C7H16O3 | 020324-32-7 | 50 |
| 3 | 2-Propanol, 1-(2-ethoxypropoxy)- | 162 | C8H18O3 | 010143-32-5 | 42 |
| 4 | Ethyl ether | 74 | C4H10O | 000060-29-7 | 38 |
| 5 | 1-Propanol, 2-(2-hydroxypropoxy)- | 134 | C6H14O3 | 000106-62-7 | 37 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

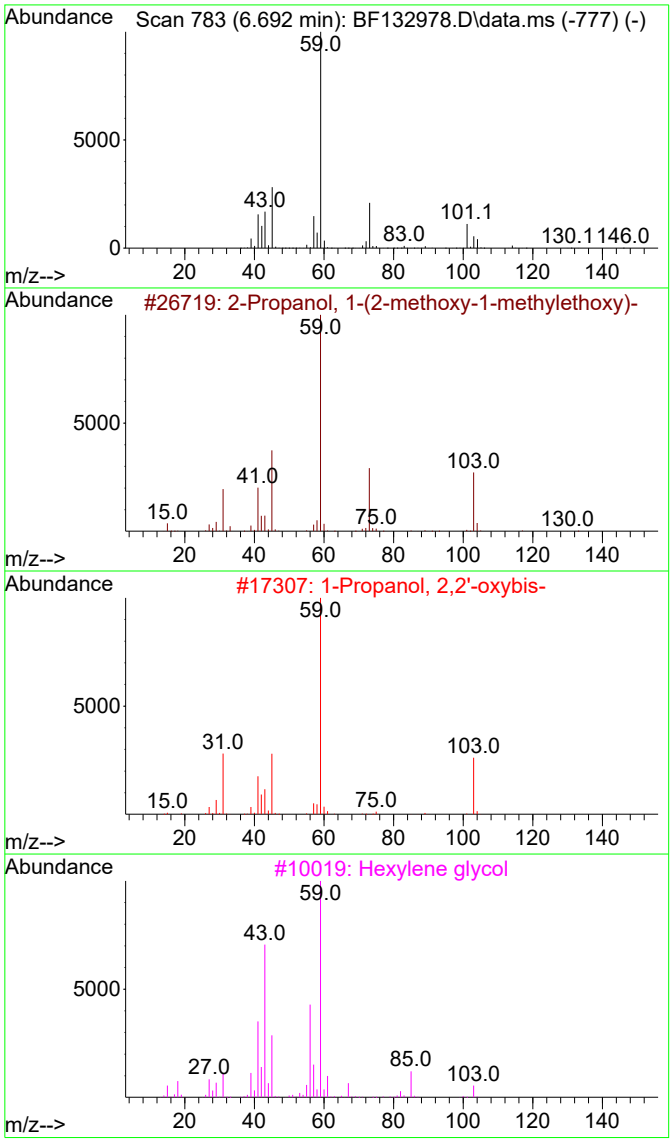
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

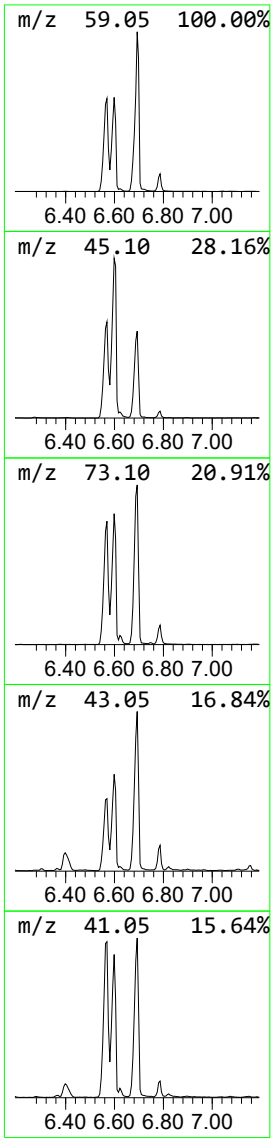
**********

Peak Number 7 1-Propanol, 2,2'-oxybis- Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.692 | 78.92 ng | 4966840 | 1,4-Dichlorobenzene-d4 | 6.751 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Propanol, 1-(2-methoxy-1-methy... | 148 | C7H16O3 | 020324-32-7 | 64 |
| 2 | 1-Propanol, 2,2'-oxybis- | 134 | C6H14O3 | 000108-61-2 | 59 |
| 3 | Hexylene glycol | 118 | C6H14O2 | 000107-41-5 | 59 |
| 4 | 1-Propanol, 2-(2-hydroxypropoxy)- | 134 | C6H14O3 | 000106-62-7 | 53 |
| 5 | (R)-(-)-2-Methyl-2,4-pentanediol | 118 | C6H14O2 | 099210-90-9 | 45 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

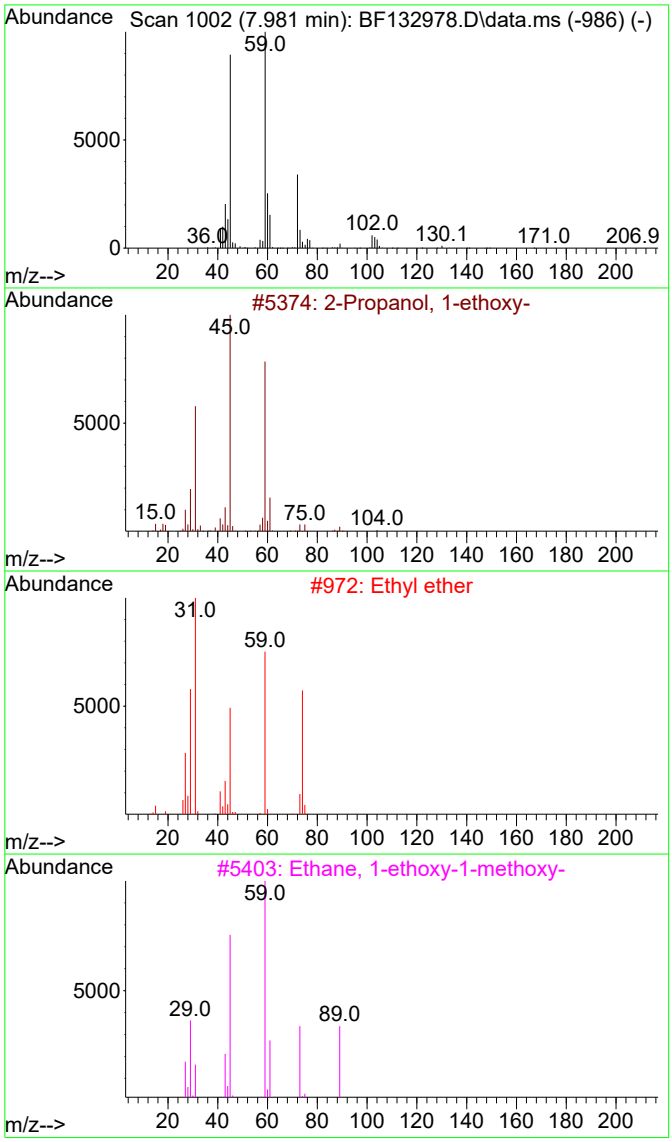
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

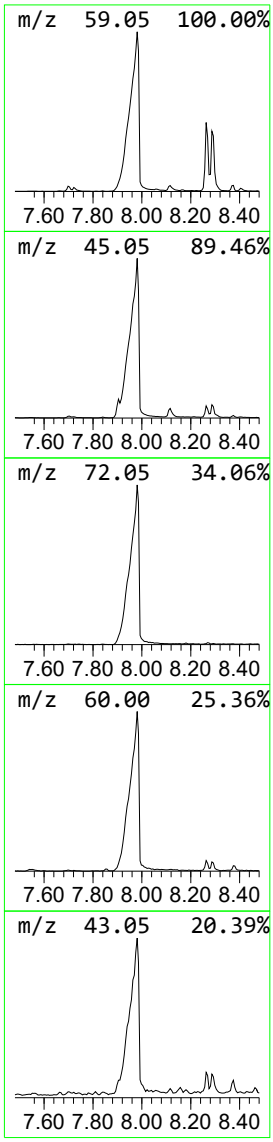
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 2-Propanol, 1-ethoxy- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.981 | 55.52 ng | 4711650 | Naphthalene-d8 | 8.034 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Propanol, 1-ethoxy- | 104 | C5H12O2 | 001569-02-4 | 59 |
| 2 | Ethyl ether | 74 | C4H10O | 000060-29-7 | 58 |
| 3 | Ethane, 1-ethoxy-1-methoxy- | 104 | C5H12O2 | 010471-14-4 | 36 |
| 4 | Ethanol, 2-[2-(2-methoxyethoxy)e... | 164 | C7H16O4 | 000112-35-6 | 36 |
| 5 | Acetic acid, ethoxy-, ethyl ester | 132 | C6H12O3 | 000817-95-8 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

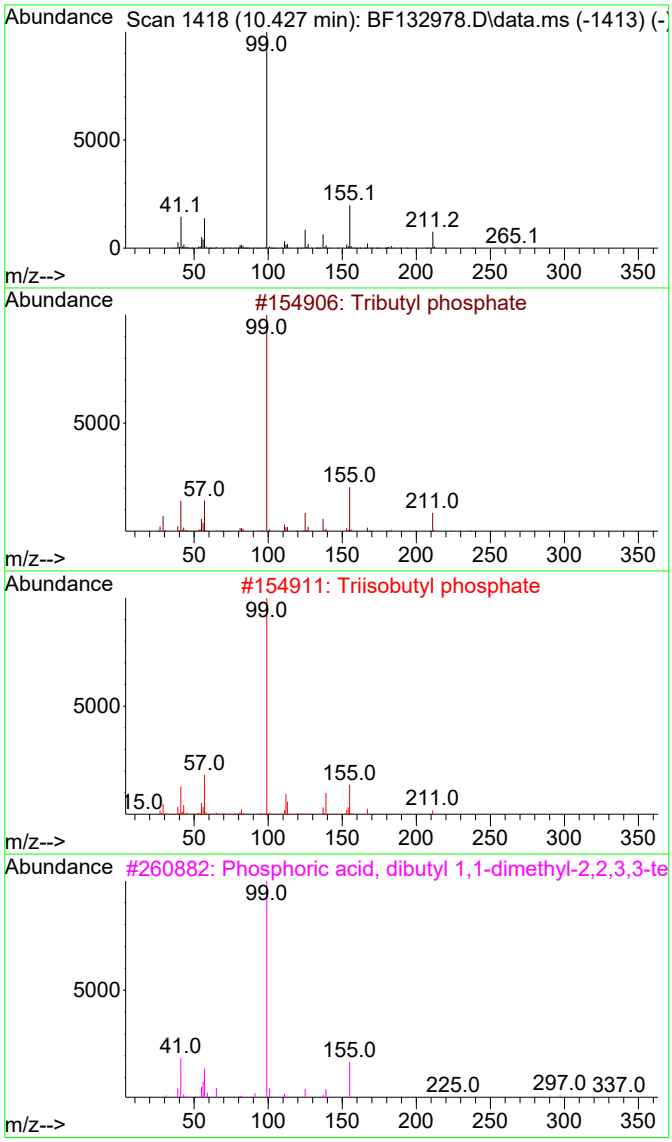
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

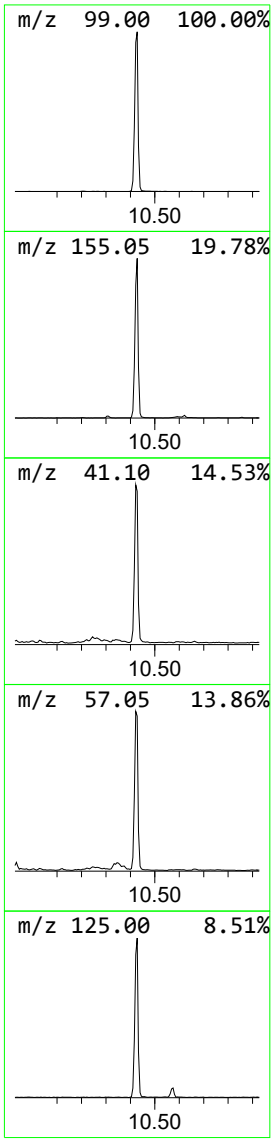
**********

Peak Number 9 Tributyl phosphate Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.428 | 37.63 ng | 3454560 | Acenaphthene-d10 | 9.786 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tributyl phosphate | 266 | C12H27O4P | 000126-73-8 | 91 |
| 2 | Triisobutyl phosphate | 266 | C12H27O4P | 000126-71-6 | 72 |
| 3 | Phosphoric acid, dibutyl 1,1-dim... | 352 | C13H25F4O4P | 1000298-93-9 | 72 |
| 4 | Phosphoric acid, dibutyl 3-trifl... | 348 | C14H28F3O4P | 1000298-93-8 | 43 |
| 5 | Butyl diethyl phosphate | 210 | C8H19O4P | 1000298-58-0 | 39 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

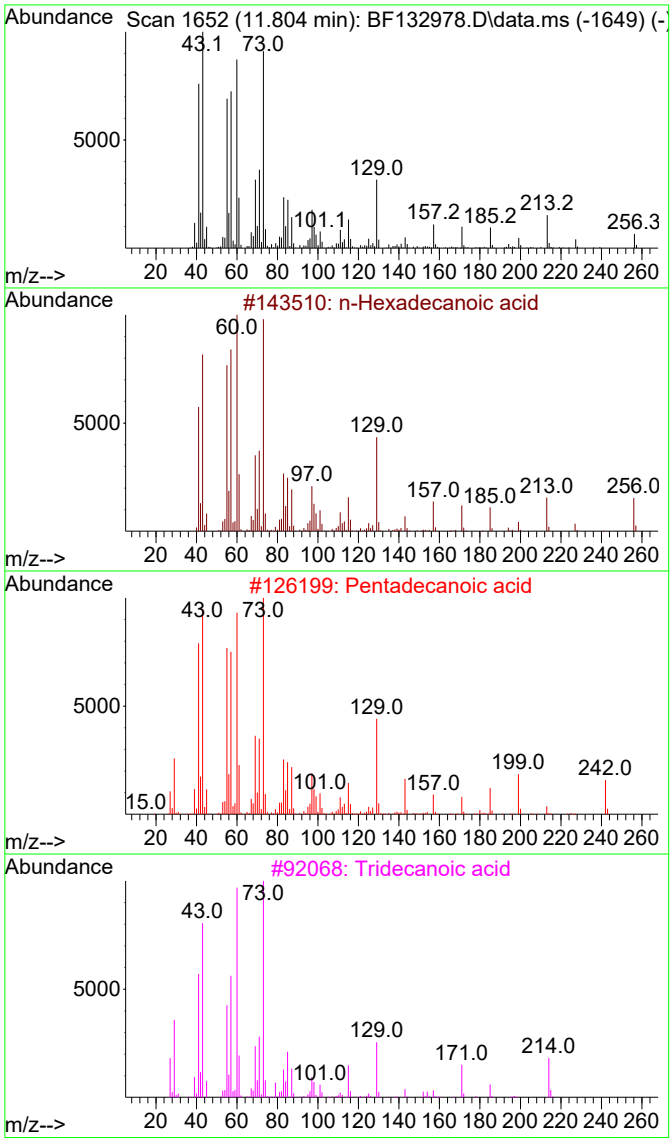
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

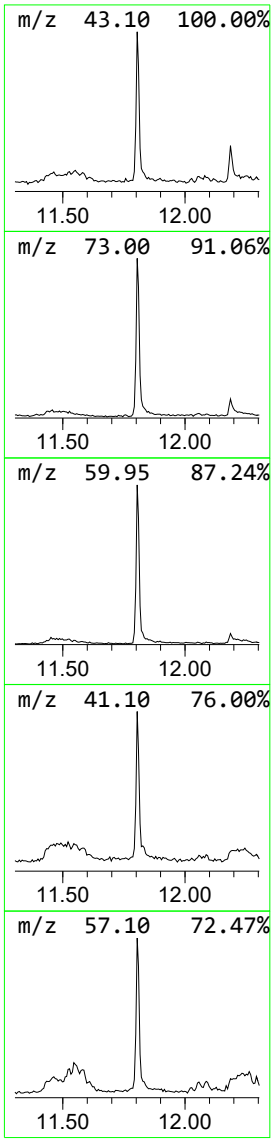
**********

Peak Number 10 n-Hexadecanoic acid Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.804 | 7.42 ng | 626852 | Phenanthrene-d10 | 11.269 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99 |
| 2 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 87 |
| 3 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 81 |
| 4 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 72 |
| 5 | Isopropyl palmitate | 298 | C19H38O2 | 000142-91-6 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

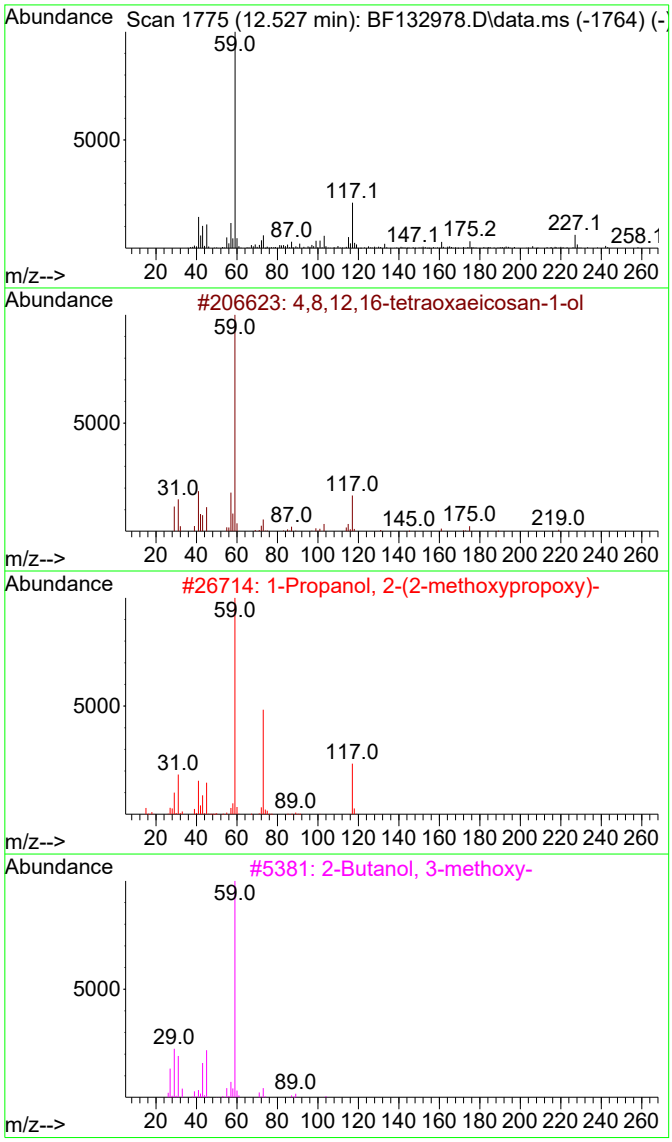
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

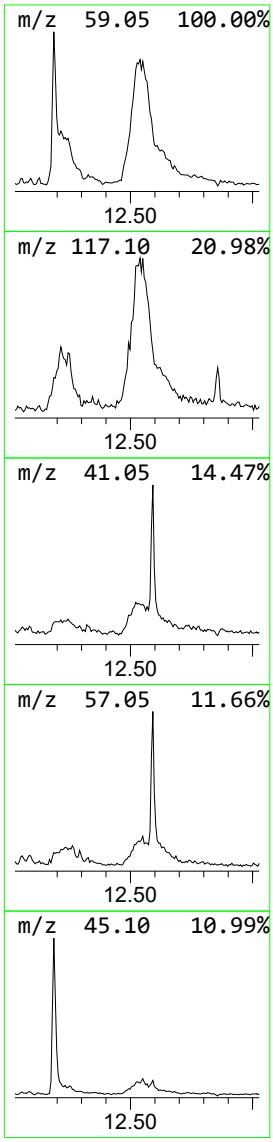
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 11 4,8,12,16-tetraoxaeicosan-1-ol Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.527 | 6.24 ng | 526969 | Phenanthrene-d10 | 11.269 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4,8,12,16-tetraoxaeicosan-1-ol | 306 | C16H34O5 | 1010397-63-6 | 72 |
| 2 | 1-Propanol, 2-(2-methoxypropoxy)- | 148 | C7H16O3 | 013588-28-8 | 59 |
| 3 | 2-Butanol, 3-methoxy- | 104 | C5H12O2 | 053778-72-6 | 50 |
| 4 | 2-Propanol, 1-[1-methyl-2-(2-pro... | 174 | C9H18O3 | 055956-25-7 | 50 |
| 5 | Butanamide, 3-methyl- | 101 | C5H11NO | 000541-46-8 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

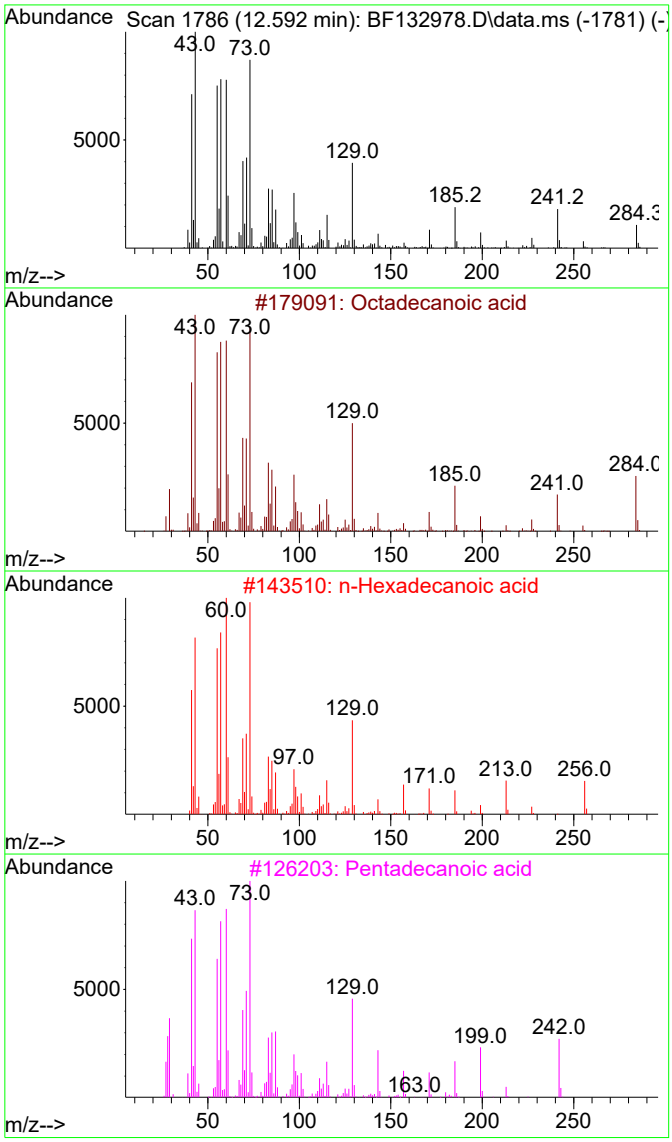
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

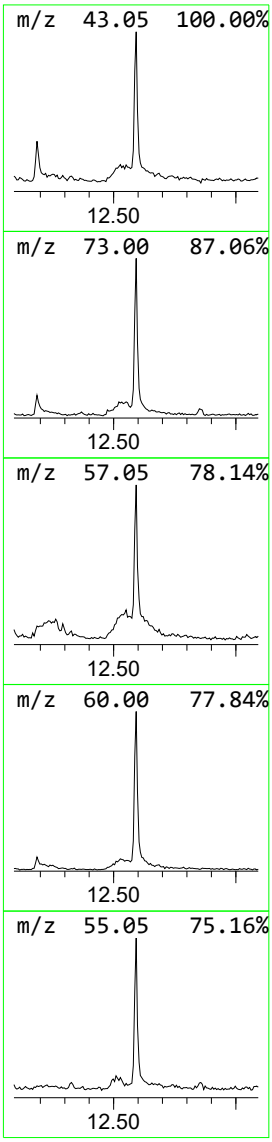
**********

Peak Number 12 Octadecanoic acid Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.592 | 14.16 ng | 787819 | Chrysene-d12 | 13.898 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 99 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 93 |
| 3 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 86 |
| 4 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 83 |
| 5 | Octadecanoic acid, 2-(2-hydroxye... | 372 | C22H44O4 | 000106-11-6 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

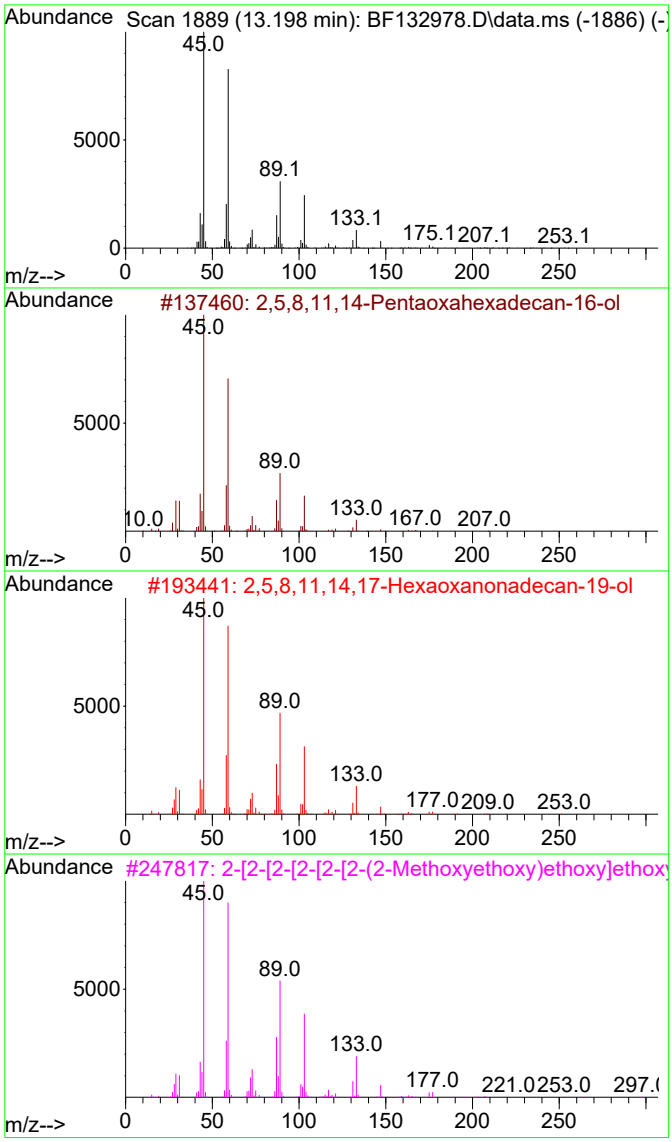
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

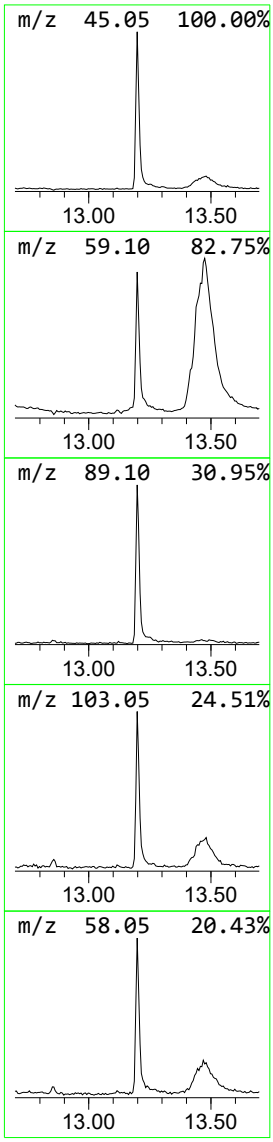
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 13 2,5,8,11,14-Pentaoxahexadec... Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.198 | 9.89 ng | 550229 | Chrysene-d12 | 13.898 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,5,8,11,14-Pentaoxahexadecan-16-ol | 252 | C11H24O6 | 023778-52-1 | 91 |
| 2 | 2,5,8,11,14,17-Hexaoxanonadecan-... | 296 | C13H28O7 | 023601-40-3 | 90 |
| 3 | 2-[2-[2-[2-[2-[2-(2-Methoxyethox... | 340 | C15H32O8 | 004437-01-8 | 90 |
| 4 | Methyl 2,5,8,11,14,17-hexaoxanon... | 324 | C14H28O8 | 1000366-81-4 | 52 |
| 5 | Methyl 2,5,8,11,14,17,20,23-octa... | 412 | C18H36O10 | 1000366-81-2 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

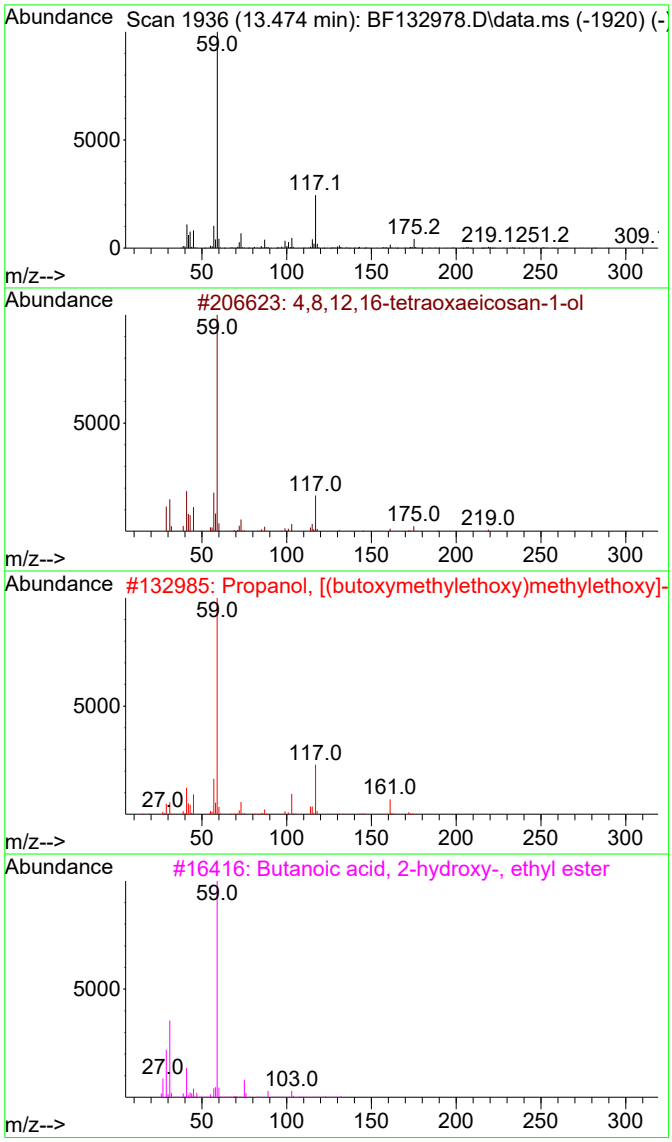
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

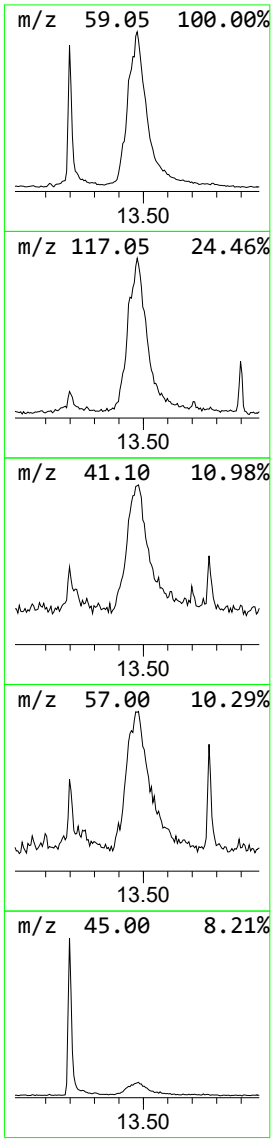
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 14 Propanol, [(butoxymethyleth... Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.474 | 28.10 ng | 1563550 | Chrysene-d12 | 13.898 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4,8,12,16-tetraoxaeicosan-1-ol | 306 | C16H34O5 | 1010397-63-6 | 64 |
| 2 | Propanol, [(butoxymethylethoxy)m... | 248 | C13H28O4 | 055934-93-5 | 56 |
| 3 | Butanoic acid, 2-hydroxy-, ethyl... | 132 | C6H12O3 | 052089-54-0 | 50 |
| 4 | Pentane, 2-methoxy- | 102 | C6H14O | 006795-88-6 | 50 |
| 5 | 2-Butanol, 3-methoxy- | 104 | C5H12O2 | 053778-72-6 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

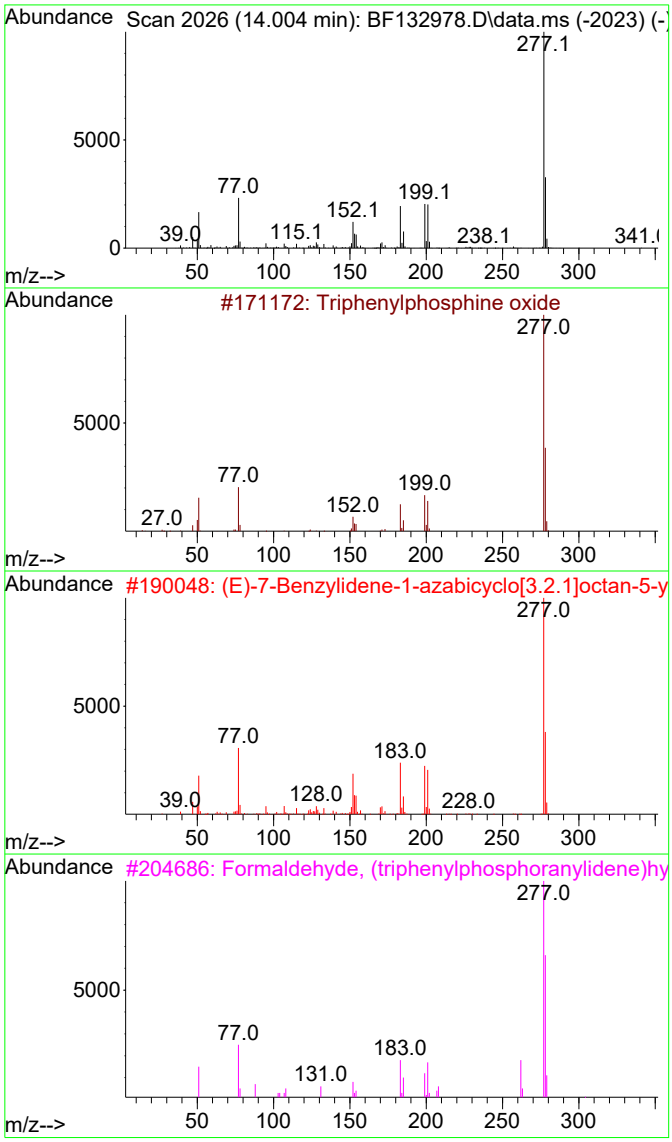
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

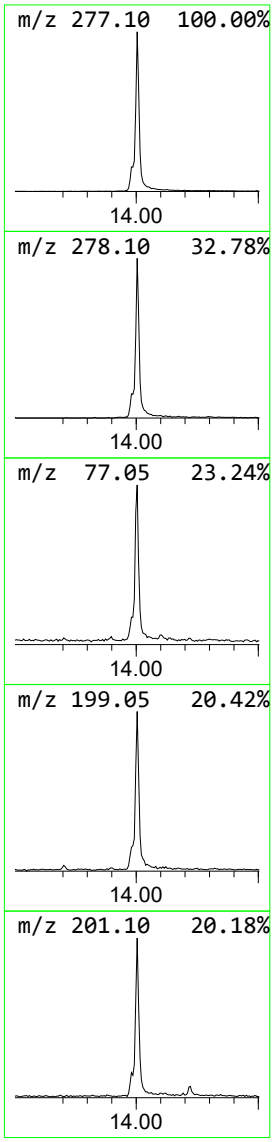
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 Triphenylphosphine oxide Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.004 | 12.18 ng | 677935 | Chrysene-d12 | 13.898 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 96 |
| 2 | (E)-7-Benzylidene-1-azabicyclo[3... | 293 | C15H19NO3S | 1000483-83-4 | 87 |
| 3 | Formaldehyde, (triphenylphosphor... | 304 | C19H17N2P | 015990-54-2 | 50 |
| 4 | Carbazol-1-one, 1,2,3,4-tetrahyd... | 277 | C18H15NO2 | 299962-12-2 | 25 |
| 5 | 4-Nitro-4'-methyldiphenylsulfone | 277 | C13H11NO4S | 004094-37-5 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

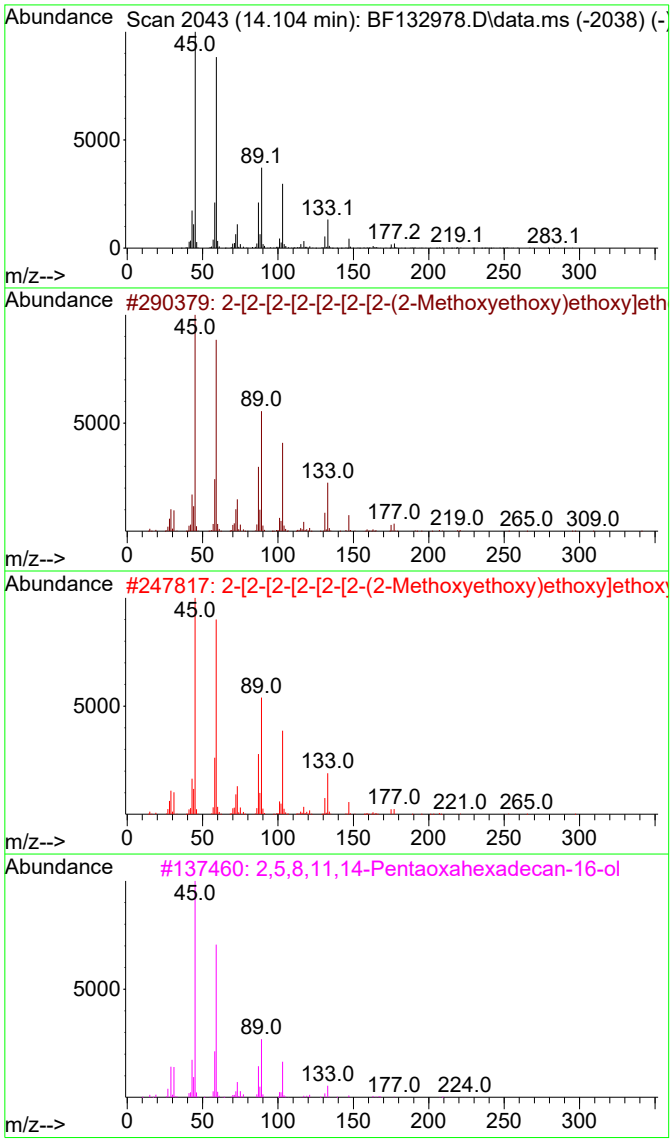
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

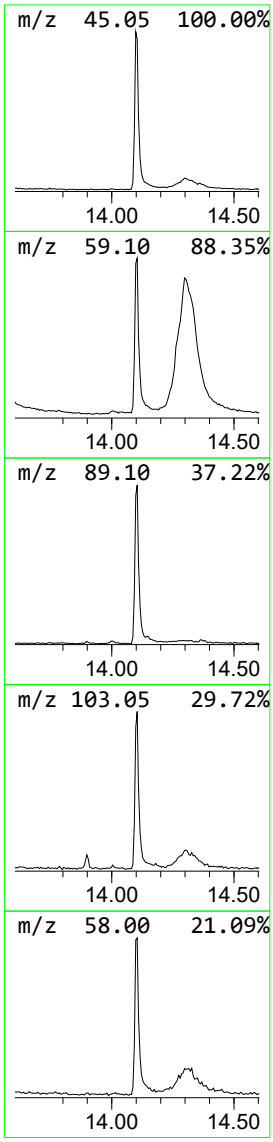
**********

Peak Number 16 unknown14.104 Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.104 | 13.92 ng | 774366 | Chrysene-d12 | 13.898 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-[2-[2-[2-[2-[2-[2-(2-Methoxyet... | 384 | C17H36O9 | 025990-96-9 | 91 |
| 2 | 2-[2-[2-[2-[2-[2-(2-Methoxyethox... | 340 | C15H32O8 | 004437-01-8 | 91 |
| 3 | 2,5,8,11,14-Pentaoxahexadecan-16-ol | 252 | C11H24O6 | 023778-52-1 | 86 |
| 4 | 2-[2-[2-[2-[2-[2-[2-[2-[2-(2-Met... | 472 | C21H44O11 | 027425-92-9 | 81 |
| 5 | 2-[2-[2-[2-[2-[2-[2-[2-[2-[2-(2-... | 516 | C23H48O12 | 1010352-04-4 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

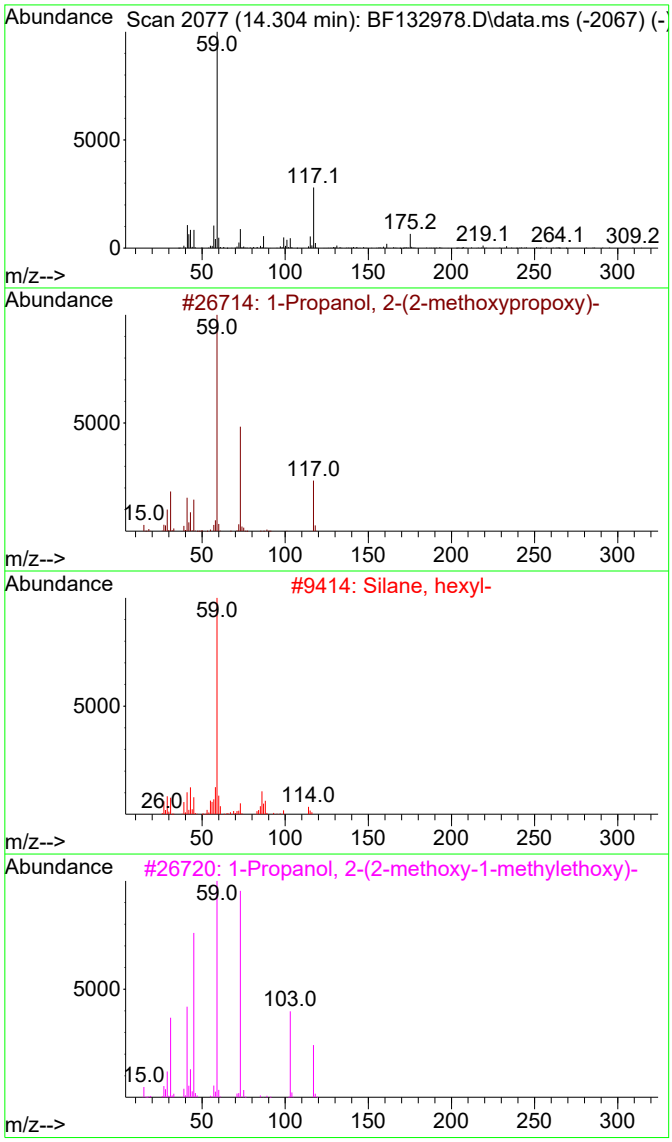
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

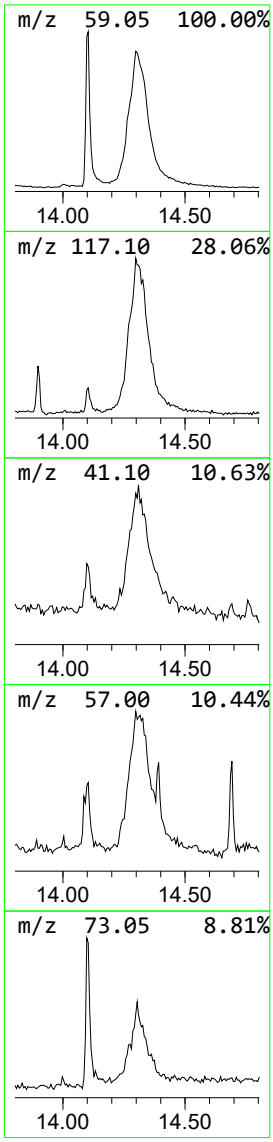
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 17 1-Propanol, 2-(2-methoxypro... Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.304 | 16.59 ng | 923061 | Chrysene-d12 | 13.898 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Propanol, 2-(2-methoxypropoxy)- | 148 | C7H16O3 | 013588-28-8 | 59 |
| 2 | Silane, hexyl- | 116 | C6H16Si | 001072-14-6 | 53 |
| 3 | 1-Propanol, 2-(2-methoxy-1-methy... | 148 | C7H16O3 | 055956-21-3 | 50 |
| 4 | 2-Propanol, 1-(2-methoxypropoxy)- | 148 | C7H16O3 | 013429-07-7 | 50 |
| 5 | 3-Hexanol | 102 | C6H14O | 000623-37-0 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

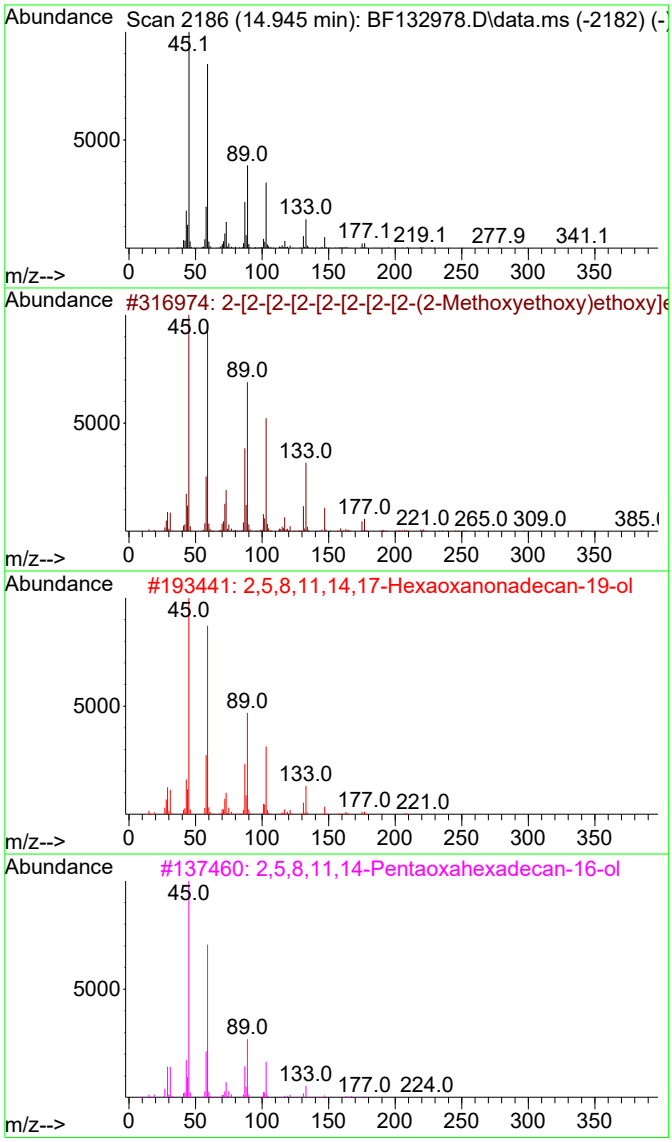
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

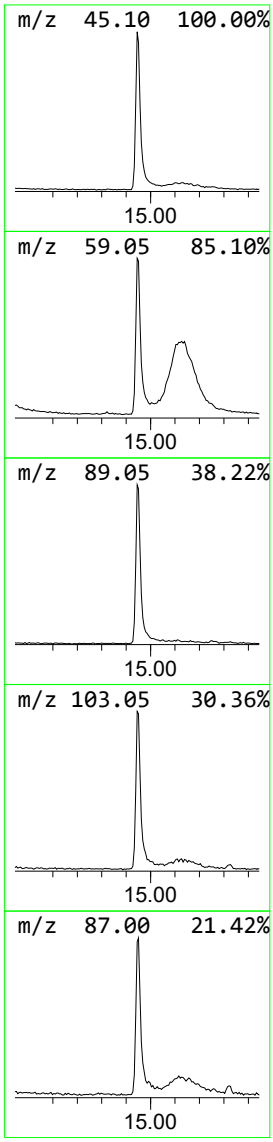
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 18 2,5,8,11,14,17-Hexaoxanonad... Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.945 | 19.65 ng | 1026740 | Perylene-d12 | 15.327 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-[2-[2-[2-[2-[2-[2-[2-(2-Methox... | 428 | C19H40O10 | 006048-68-6 | 91 |
| 2 | 2,5,8,11,14,17-Hexaoxanonadecan-... | 296 | C13H28O7 | 023601-40-3 | 78 |
| 3 | 2,5,8,11,14-Pentaoxahexadecan-16-ol | 252 | C11H24O6 | 023778-52-1 | 72 |
| 4 | 2-[2-[2-[2-[2-[2-[2-[2-[2-(2-Met... | 472 | C21H44O11 | 027425-92-9 | 70 |
| 5 | 2-[2-[2-[2-[2-(2-Methoxyethox... | 340 | C15H32O8 | 004437-01-8 | 70 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

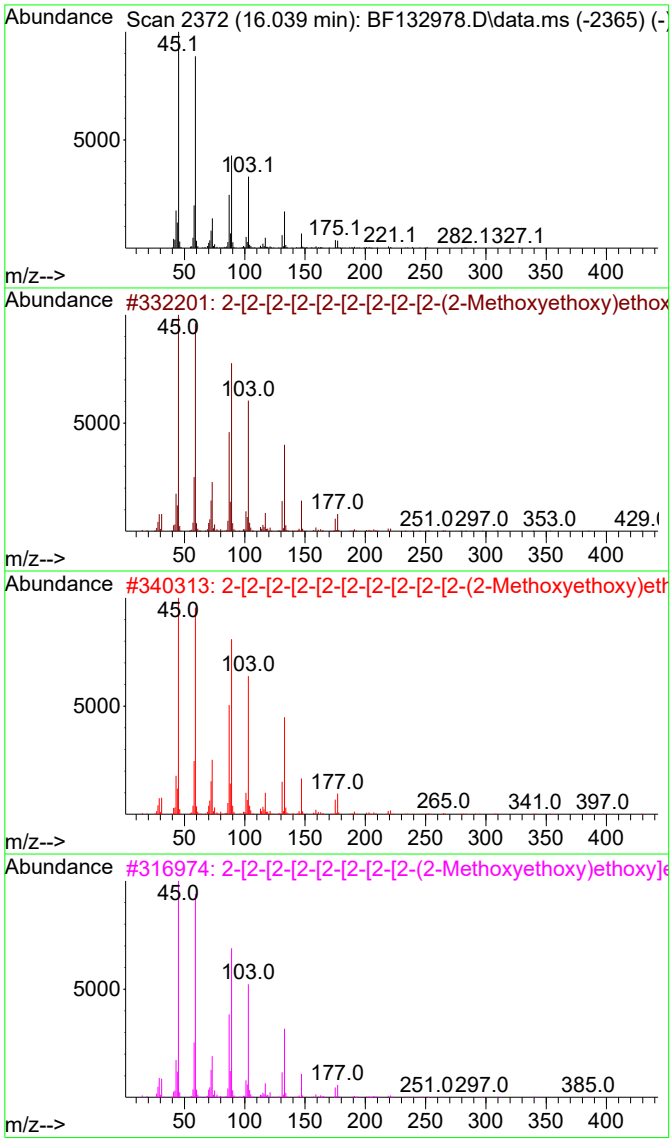
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

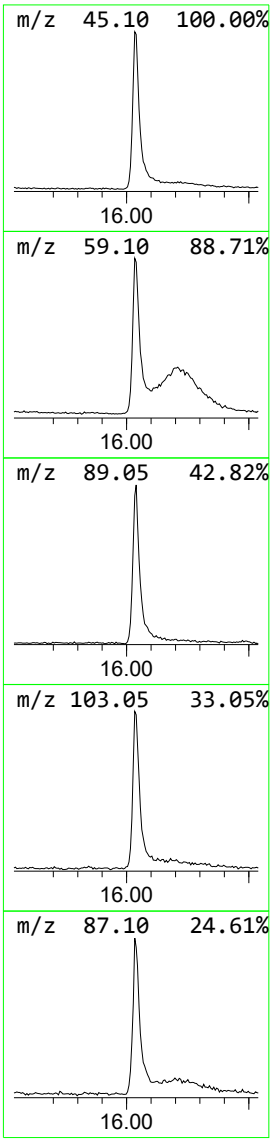
**********

Peak Number 19 unknown16.039 Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.039 | 16.94 ng | 885244 | Perylene-d12 | 15.327 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-[2-[2-[2-[2-[2-[2-[2-[2-(2-Met... | 472 | C21H44O11 | 027425-92-9 | 91 |
| 2 | 2-[2-[2-[2-[2-[2-[2-[2-[2-[2-(2-... | 516 | C23H48O12 | 1010352-04-4 | 91 |
| 3 | 2-[2-[2-[2-[2-[2-[2-[2-(2-Methox... | 428 | C19H40O10 | 006048-68-6 | 91 |
| 4 | 2,5,8,11,14-Pentaoxahexadecan-16-ol | 252 | C11H24O6 | 023778-52-1 | 86 |
| 5 | 2,5,8,11,14,17-Hexaoxanonadecan-... | 296 | C13H28O7 | 023601-40-3 | 78 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

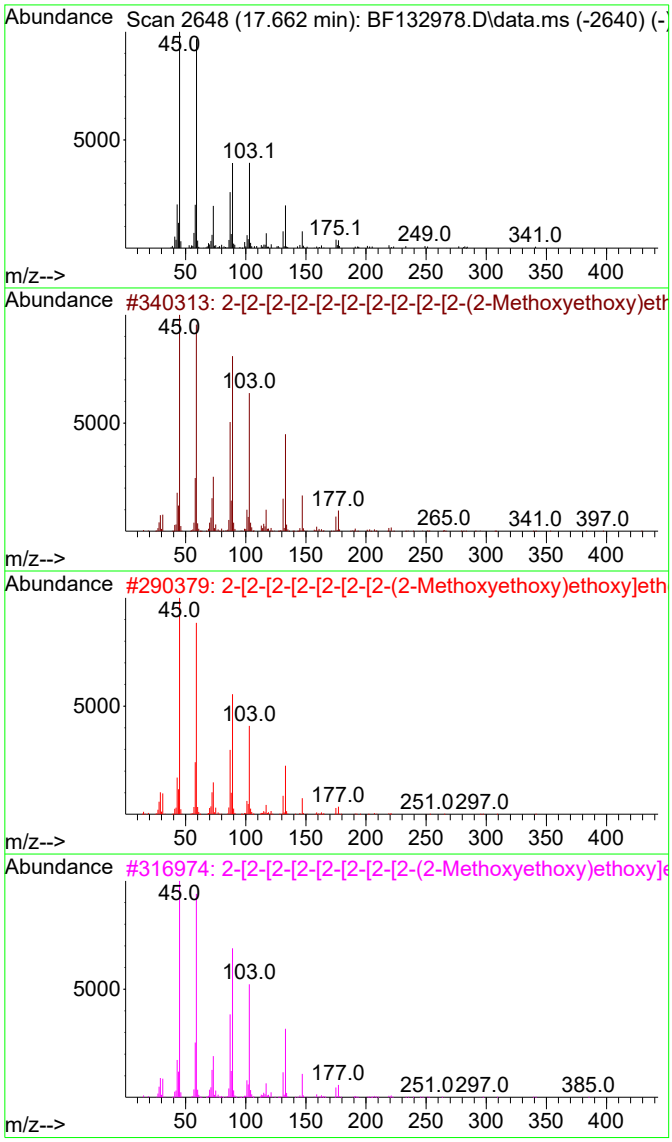
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

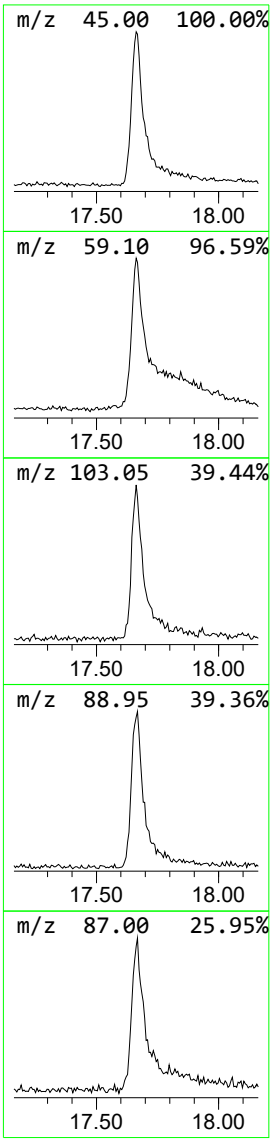
**********

Peak Number 20 unknown17.662 Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.662 | 8.93 ng | 466456 | Perylene-d12 | 15.327 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-[2-[2-[2-[2-[2-[2-[2-[2-[2-(2-... | 516 | C23H48O12 | 1010352-04-4 | 91 |
| 2 | 2-[2-[2-[2-[2-[2-[2-(2-Methoxyet... | 384 | C17H36O9 | 025990-96-9 | 72 |
| 3 | 2-[2-[2-[2-[2-[2-[2-[2-(2-Methox... | 428 | C19H40O10 | 006048-68-6 | 62 |
| 4 | 2,5,8,11,14-Pentaoxahexadecan-16-ol | 252 | C11H24O6 | 023778-52-1 | 53 |
| 5 | 2-[2-[2-[2-[2-[2-[2-[2-[2-(2-Met... | 472 | C21H44O11 | 027425-92-9 | 53 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042123\
Data File : BF132978.D
Acq On : 21 Apr 2023 20:01
Operator : CG\JU
Sample : 02415-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
SAMPLE-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.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-ethoxy- | 2.540 | 18.8 | ng | 1186070 | 1 | 6.751 | 1258780 | 20.0 |
| 2-Pentanone, 4-... | 4.945 | 25.9 | ng | 1631890 | 1 | 6.751 | 1258780 | 20.0 |
| o-Xylene | 5.592 | 5.7 | ng | 359409 | 1 | 6.751 | 1258780 | 20.0 |
| 2-Propanol, 1-b... | 6.051 | 7.4 | ng | 468346 | 1 | 6.751 | 1258780 | 20.0 |
| 2-Propanol, 1-(... | 6.569 | 61.1 | ng | 3845410 | 1 | 6.751 | 1258780 | 20.0 |
| Ethane, 1-ethox... | 6.598 | 60.0 | ng | 3776180 | 1 | 6.751 | 1258780 | 20.0 |
| 1-Propanol, 2,2... | 6.692 | 78.9 | ng | 4966840 | 1 | 6.751 | 1258780 | 20.0 |
| 2-Propanol, 1-e... | 7.981 | 55.5 | ng | 4711650 | 2 | 8.034 | 1697320 | 20.0 |
| Tributyl phosphate | 10.428 | 37.6 | ng | 3454560 | 3 | 9.786 | 1836020 | 20.0 |
| n-Hexadecanoic ... | 11.804 | 7.4 | ng | 626852 | 4 | 11.269 | 1688970 | 20.0 |
| 4,8,12,16-tetra... | 12.527 | 6.2 | ng | 526969 | 4 | 11.269 | 1688970 | 20.0 |
| Octadecanoic acid | 12.592 | 14.2 | ng | 787819 | 5 | 13.898 | 1112800 | 20.0 |
| 2,5,8,11,14-Pen... | 13.198 | 9.9 | ng | 550229 | 5 | 13.898 | 1112800 | 20.0 |
| Propanol, [(but... | 13.474 | 28.1 | ng | 1563550 | 5 | 13.898 | 1112800 | 20.0 |
| Triphenylphosph... | 14.004 | 12.2 | ng | 677935 | 5 | 13.898 | 1112800 | 20.0 |
| unknown14.104 | 14.104 | 13.9 | ng | 774366 | 5 | 13.898 | 1112800 | 20.0 |
| 1-Propanol, 2-(... | 14.304 | 16.6 | ng | 923061 | 5 | 13.898 | 1112800 | 20.0 |
| 2,5,8,11,14,17-... | 14.945 | 19.6 | ng | 1026740 | 6 | 15.327 | 1045160 | 20.0 |
| unknown16.039 | 16.039 | 16.9 | ng | 885244 | 6 | 15.327 | 1045160 | 20.0 |
| unknown17.662 | 17.662 | 8.9 | ng | 466456 | 6 | 15.327 | 1045160 | 20.0 |