PDB file of KAP8.1 Model

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REMARK   Accelrys Discovery Studio PDB file
REMARK   Created:  2016-07-01T15:14:01Z
SSBOND   1 CYS A    1    CYS A   28
SSBOND   2 CYS A   13    CYS A   43
ATOM      1  N   MET A  -1      27.440  17.934  11.835  1.00  0.00
N1+
ATOM      2  HT1 MET A  -1      27.384  18.772  12.446  1.00  0.00
H
ATOM      3  HT2 MET A  -1      28.365  17.475  11.936  1.00  0.00
H
ATOM      4  HT3 MET A  -1      26.717  17.236  12.139  1.00  0.00
H
ATOM      5  CA  MET A  -1      27.128  18.212  10.432  1.00  0.00
C
ATOM      6  HA  MET A  -1      26.559  19.143  10.391  1.00  0.00
H
ATOM      7  CB  MET A  -1      28.350  18.270   9.493  1.00  0.00
C
ATOM      8  HB1 MET A  -1      29.077  17.497   9.747  1.00  0.00
H
ATOM      9  HB2 MET A  -1      28.041  18.058   8.467  1.00  0.00
H
ATOM     10  CG  MET A  -1      29.020  19.648   9.471  1.00  0.00
C
ATOM     11  HG1 MET A  -1      28.276  20.420   9.270  1.00  0.00
H
ATOM     12  HG2 MET A  -1      29.478  19.862  10.437  1.00  0.00
H
ATOM     13  SD  MET A  -1      30.292  19.698   8.194  1.00  0.00
S
ATOM     14  CE  MET A  -1      29.813  21.223   7.360  1.00  0.00
C
ATOM     15  HE1 MET A  -1      30.570  21.513   6.630  1.00  0.00
H
ATOM     16  HE2 MET A  -1      29.713  22.029   8.088  1.00  0.00
H
ATOM     17  HE3 MET A  -1      28.848  21.100   6.865  1.00  0.00
H
ATOM     18  C   MET A  -1      26.167  17.100  10.050  1.00  0.00
C
ATOM     19  O   MET A  -1      25.973  16.228  10.887  1.00  0.00
O
ATOM     20  N   LEU A   0      25.577  17.246   8.844  1.00  0.00
N
ATOM     21  HN  LEU A   0      25.973  17.949   8.252  1.00  0.00
H
ATOM     22  CA  LEU A   0      24.200  16.840   8.482  1.00  0.00
C
ATOM     23  HA  LEU A   0      24.212  16.663   7.407  1.00  0.00
H
ATOM     24  CB  LEU A   0      23.629  15.615   9.238  1.00  0.00
C
ATOM     25  HB1 LEU A   0      23.763  15.734  10.315  1.00  0.00
H
ATOM     26  HB2 LEU A   0      22.549  15.557   9.112  1.00  0.00
H
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ATOM     27  CG  LEU A   0      24.188  14.265   8.765  1.00  0.00
C
ATOM     28  HG  LEU A   0      25.269  14.352   8.652  1.00  0.00
H
ATOM     29  CD1 LEU A   0      23.932  13.186   9.822  1.00  0.00
C
ATOM     30 1HD1 LEU A   0      24.258  12.205   9.478  1.00  0.00
H
ATOM     31 2HD1 LEU A   0      22.875  13.108  10.074  1.00  0.00
H
ATOM     32 3HD1 LEU A   0      24.471  13.411  10.742  1.00  0.00
H
ATOM     33  CD2 LEU A   0      23.576  13.841   7.420  1.00  0.00
C
ATOM     34 1HD2 LEU A   0      23.985  12.880   7.101  1.00  0.00
H
ATOM     35 2HD2 LEU A   0      23.789  14.562   6.631  1.00  0.00
H
ATOM     36 3HD2 LEU A   0      22.491  13.728   7.488  1.00  0.00
H
ATOM     37  C   LEU A   0      23.374  18.109   8.728  1.00  0.00
C
ATOM     38  O   LEU A   0      23.998  19.155   8.908  1.00  0.00
O
ATOM     39  N   CYS A   1      22.034  17.996   8.789  1.00  0.00
N
ATOM     40  HN  CYS A   1      21.538  17.167   8.525  1.00  0.00
H
ATOM     41  CA  CYS A   1      21.277  19.061   9.420  1.00  0.00
C
ATOM     42  HA  CYS A   1      21.905  19.542  10.176  1.00  0.00
H
ATOM     43  CB  CYS A   1      20.820  20.113   8.404  1.00  0.00
C
ATOM     44  HB1 CYS A   1      19.777  20.387   8.574  1.00  0.00
H
ATOM     45  HB2 CYS A   1      21.406  21.022   8.535  1.00  0.00
H
ATOM     46  SG  CYS A   1      21.021  19.567   6.699  1.00  0.00
S
ATOM     47  C   CYS A   1      20.138  18.363  10.144  1.00  0.00
C
ATOM     48  O   CYS A   1      19.100  18.000   9.578  1.00  0.00
O
ATOM     49  N   ASP A   2      20.428  18.155  11.438  1.00  0.00
N
ATOM     50  HN  ASP A   2      21.322  18.406  11.821  1.00  0.00
H
ATOM     51  CA  ASP A   2      19.490  17.554  12.374  1.00  0.00
C
ATOM     52  HA  ASP A   2      19.068  16.678  11.877  1.00  0.00
H
ATOM     53  CB  ASP A   2      20.220  17.098  13.659  1.00  0.00
C
ATOM     54  HB1 ASP A   2      20.035  17.808  14.469  1.00  0.00
H
ATOM     55  HB2 ASP A   2      19.805  16.150  14.006  1.00  0.00
H
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ATOM     56  CG  ASP A   2      21.745  16.973  13.484  1.00  0.00
C
ATOM     57  OD1 ASP A   2      22.263  15.898  13.203  1.00  0.00
O
ATOM     58  OD2 ASP A   2      22.430  17.980  13.630  1.00  0.00
O1-
ATOM     59  C   ASP A   2      18.396  18.609  12.637  1.00  0.00
C
ATOM     60  O   ASP A   2      18.644  19.677  13.216  1.00  0.00
O
ATOM     61  N   ASN A   3      17.240  18.244  12.045  1.00  0.00
N
ATOM     62  HN  ASN A   3      17.105  17.275  11.848  1.00  0.00
H
ATOM     63  CA  ASN A   3      16.296  19.194  11.500  1.00  0.00
C
ATOM     64  HA  ASN A   3      16.857  20.111  11.296  1.00  0.00
H
ATOM     65  CB  ASN A   3      15.719  18.639  10.192  1.00  0.00
C
ATOM     66  HB1 ASN A   3      16.508  18.415   9.477  1.00  0.00
H
ATOM     67  HB2 ASN A   3      15.186  17.708  10.373  1.00  0.00
H
ATOM     68  CG  ASN A   3      14.775  19.662   9.557  1.00  0.00
C
ATOM     69  OD1 ASN A   3      15.197  20.733   9.106  1.00  0.00
O
ATOM     70  ND2 ASN A   3      13.472  19.316   9.601  1.00  0.00
N
ATOM     71 1HD2 ASN A   3      13.081  19.990   8.974  1.00  0.00
H
ATOM     72 2HD2 ASN A   3      12.954  18.623  10.116  1.00  0.00
H
ATOM     73  C   ASN A   3      15.156  19.498  12.470  1.00  0.00
C
ATOM     74  O   ASN A   3      14.543  18.652  13.128  1.00  0.00
O
ATOM     75  N   PHE A   4      14.880  20.824  12.472  1.00  0.00
N
ATOM     76  HN  PHE A   4      15.431  21.352  11.822  1.00  0.00
H
ATOM     77  CA  PHE A   4      13.803  21.472  13.215  1.00  0.00
C
ATOM     78  HA  PHE A   4      14.282  21.919  14.082  1.00  0.00
H
ATOM     79  CB  PHE A   4      13.226  22.614  12.369  1.00  0.00
C
ATOM     80  HB1 PHE A   4      12.717  22.196  11.500  1.00  0.00
H
ATOM     81  HB2 PHE A   4      12.488  23.185  12.934  1.00  0.00
H
ATOM     82  CG  PHE A   4      14.325  23.534  11.912  1.00  0.00
C
ATOM     83  CD1 PHE A   4      14.726  23.530  10.581  1.00  0.00
C
ATOM     84  HD1 PHE A   4      14.248  22.885   9.859  1.00  0.00
H
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ATOM     85  CD2 PHE A   4      14.977  24.374  12.812  1.00  0.00
C
ATOM     86  HD2 PHE A   4      14.684  24.403  13.855  1.00  0.00
H
ATOM     87  CE1 PHE A   4      15.761  24.342  10.148  1.00  0.00
C
ATOM     88  HE1 PHE A   4      16.056  24.292   9.110  1.00  0.00
H
ATOM     89  CE2 PHE A   4      16.021  25.181  12.380  1.00  0.00
C
ATOM     90  HE2 PHE A   4      16.549  25.812  13.081  1.00  0.00
H
ATOM     91  CZ  PHE A   4      16.412  25.168  11.048  1.00  0.00
C
ATOM     92  HZ  PHE A   4      17.226  25.795  10.712  1.00  0.00
H
ATOM     93  C   PHE A   4      12.681  20.502  13.660  1.00  0.00
C
ATOM     94  O   PHE A   4      11.723  20.229  12.929  1.00  0.00
O
ATOM     95  N   PRO A   5      12.786  20.032  14.948  1.00  0.00
N
ATOM     96  CD  PRO A   5      13.933  20.257  15.835  1.00  0.00
C
ATOM     97  HD1 PRO A   5      14.857  19.972  15.327  1.00  0.00
H
ATOM     98  HD2 PRO A   5      14.006  21.303  16.133  1.00  0.00
H
ATOM     99  CA  PRO A   5      11.805  19.105  15.499  1.00  0.00
C
ATOM    100  HA  PRO A   5      11.466  18.424  14.722  1.00  0.00
H
ATOM    101  CB  PRO A   5      12.651  18.334  16.528  1.00  0.00
C
ATOM    102  HB1 PRO A   5      12.103  17.822  17.317  1.00  0.00
H
ATOM    103  HB2 PRO A   5      13.202  17.557  15.999  1.00  0.00
H
ATOM    104  CG  PRO A   5      13.671  19.354  17.041  1.00  0.00
C
ATOM    105  HG1 PRO A   5      13.236  19.944  17.850  1.00  0.00
H
ATOM    106  HG2 PRO A   5      14.584  18.890  17.420  1.00  0.00
H
ATOM    107  C   PRO A   5      10.545  19.811  16.082  1.00  0.00
C
ATOM    108  O   PRO A   5       9.832  20.550  15.403  1.00  0.00
O
ATOM    109  N   GLY A   6      10.233  19.514  17.358  1.00  0.00
N
ATOM    110  HN  GLY A   6      10.853  18.946  17.900  1.00  0.00
H
ATOM    111  CA  GLY A   6       8.891  19.834  17.850  1.00  0.00
C
ATOM    112  HA1 GLY A   6       8.197  19.942  17.014  1.00  0.00
H
ATOM    113  HA2 GLY A   6       8.563  18.989  18.447  1.00  0.00
H
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ATOM    114  C   GLY A   6       8.875  21.088  18.717  1.00  0.00
C
ATOM    115  O   GLY A   6       9.101  21.051  19.923  1.00  0.00
O
ATOM    116  N   ALA A   7       8.575  22.232  18.073  1.00  0.00
N
ATOM    117  HN  ALA A   7       8.385  22.171  17.095  1.00  0.00
H
ATOM    118  CA  ALA A   7       8.583  23.503  18.805  1.00  0.00
C
ATOM    119  HA  ALA A   7       9.590  23.604  19.215  1.00  0.00
H
ATOM    120  CB  ALA A   7       8.298  24.674  17.849  1.00  0.00
C
ATOM    121  HB1 ALA A   7       9.024  24.689  17.033  1.00  0.00
H
ATOM    122  HB2 ALA A   7       8.341  25.637  18.366  1.00  0.00
H
ATOM    123  HB3 ALA A   7       7.303  24.571  17.412  1.00  0.00
H
ATOM    124  C   ALA A   7       7.577  23.431  19.976  1.00  0.00
C
ATOM    125  O   ALA A   7       6.544  22.748  19.903  1.00  0.00
O
ATOM    126  N   VAL A   8       7.961  24.101  21.094  1.00  0.00
N
ATOM    127  HN  VAL A   8       8.754  24.719  21.072  1.00  0.00
H
ATOM    128  CA  VAL A   8       7.206  23.887  22.318  1.00  0.00
C
ATOM    129  HA  VAL A   8       7.079  22.807  22.427  1.00  0.00
H
ATOM    130  CB  VAL A   8       8.006  24.412  23.531  1.00  0.00
C
ATOM    131  HB  VAL A   8       7.410  24.208  24.421  1.00  0.00
H
ATOM    132  CG1 VAL A   8       9.322  23.637  23.699  1.00  0.00
C
ATOM    133 1HG1 VAL A   8       9.955  23.680  22.810  1.00  0.00
H
ATOM    134 2HG1 VAL A   8       9.114  22.594  23.932  1.00  0.00
H
ATOM    135 3HG1 VAL A   8       9.902  24.037  24.531  1.00  0.00
H
ATOM    136  CG2 VAL A   8       8.267  25.931  23.506  1.00  0.00
C
ATOM    137 1HG2 VAL A   8       8.843  26.240  24.380  1.00  0.00
H
ATOM    138 2HG2 VAL A   8       7.336  26.497  23.520  1.00  0.00
H
ATOM    139 3HG2 VAL A   8       8.828  26.226  22.620  1.00  0.00
H
ATOM    140  C   VAL A   8       5.814  24.536  22.214  1.00  0.00
C
ATOM    141  O   VAL A   8       5.677  25.727  21.948  1.00  0.00
O
ATOM    142  N   PHE A   9       4.813  23.701  22.515  1.00  0.00
N
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ATOM    143  HN  PHE A   9       5.071  22.758  22.715  1.00  0.00
H
ATOM    144  CA  PHE A   9       3.551  24.207  23.028  1.00  0.00
C
ATOM    145  HA  PHE A   9       3.334  25.177  22.576  1.00  0.00
H
ATOM    146  CB  PHE A   9       2.480  23.130  22.726  1.00  0.00
C
ATOM    147  HB1 PHE A   9       2.618  22.753  21.713  1.00  0.00
H
ATOM    148  HB2 PHE A   9       2.602  22.292  23.402  1.00  0.00
H
ATOM    149  CG  PHE A   9       1.052  23.569  22.871  1.00  0.00
C
ATOM    150  CD1 PHE A   9       0.333  24.021  21.771  1.00  0.00
C
ATOM    151  HD1 PHE A   9       0.799  24.079  20.794  1.00  0.00
H
ATOM    152  CD2 PHE A   9       0.423  23.511  24.111  1.00  0.00
C
ATOM    153  HD2 PHE A   9       0.950  23.103  24.962  1.00  0.00
H
ATOM    154  CE1 PHE A   9      -0.992  24.410  21.919  1.00  0.00
C
ATOM    155  HE1 PHE A   9      -1.550  24.773  21.067  1.00  0.00
H
ATOM    156  CE2 PHE A   9      -0.890  23.932  24.265  1.00  0.00
C
ATOM    157  HE2 PHE A   9      -1.373  23.913  25.235  1.00  0.00
H
ATOM    158  CZ  PHE A   9      -1.600  24.374  23.165  1.00  0.00
C
ATOM    159  HZ  PHE A   9      -2.623  24.705  23.291  1.00  0.00
H
ATOM    160  C   PHE A   9       3.823  24.307  24.533  1.00  0.00
C
ATOM    161  O   PHE A   9       4.635  23.537  25.050  1.00  0.00
O
ATOM    162  N   PRO A  10       3.122  25.215  25.273  1.00  0.00
N
ATOM    163  CD  PRO A  10       2.287  26.304  24.759  1.00  0.00
C
ATOM    164  HD1 PRO A  10       2.933  27.077  24.345  1.00  0.00
H
ATOM    165  HD2 PRO A  10       1.595  25.986  23.981  1.00  0.00
H
ATOM    166  CA  PRO A  10       3.160  25.146  26.736  1.00  0.00
C
ATOM    167  HA  PRO A  10       4.186  25.075  27.104  1.00  0.00
H
ATOM    168  CB  PRO A  10       2.510  26.482  27.152  1.00  0.00
C
ATOM    169  HB1 PRO A  10       2.046  26.527  28.138  1.00  0.00
H
ATOM    170  HB2 PRO A  10       3.297  27.238  27.169  1.00  0.00
H
ATOM    171  CG  PRO A  10       1.558  26.823  26.005  1.00  0.00
C
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ATOM    172  HG1 PRO A  10       0.617  26.282  26.126  1.00  0.00
H
ATOM    173  HG2 PRO A  10       1.332  27.890  25.952  1.00  0.00
H
ATOM    174  C   PRO A  10       2.374  23.887  27.181  1.00  0.00
C
ATOM    175  O   PRO A  10       2.552  22.801  26.628  1.00  0.00
O
ATOM    176  N   GLY A  11       1.452  24.097  28.157  1.00  0.00
N
ATOM    177  HN  GLY A  11       1.267  24.988  28.581  1.00  0.00
H
ATOM    178  CA  GLY A  11       0.486  23.089  28.535  1.00  0.00
C
ATOM    179  HA1 GLY A  11       0.755  22.113  28.120  1.00  0.00
H
ATOM    180  HA2 GLY A  11       0.481  23.032  29.624  1.00  0.00
H
ATOM    181  C   GLY A  11      -0.872  23.556  28.023  1.00  0.00
C
ATOM    182  O   GLY A  11      -1.036  24.669  27.518  1.00  0.00
O
ATOM    183  N   GLY A  12      -1.810  22.617  28.188  1.00  0.00
N
ATOM    184  HN  GLY A  12      -1.746  21.952  28.941  1.00  0.00
H
ATOM    185  CA  GLY A  12      -3.132  22.787  27.625  1.00  0.00
C
ATOM    186  HA1 GLY A  12      -3.334  23.838  27.413  1.00  0.00
H
ATOM    187  HA2 GLY A  12      -3.205  22.188  26.718  1.00  0.00
H
ATOM    188  C   GLY A  12      -4.050  22.273  28.701  1.00  0.00
C
ATOM    189  O   GLY A  12      -3.674  22.228  29.867  1.00  0.00
O
ATOM    190  N   CYS A  13      -5.262  21.896  28.321  1.00  0.00
N
ATOM    191  HN  CYS A  13      -5.544  22.011  27.368  1.00  0.00
H
ATOM    192  CA  CYS A  13      -6.050  21.012  29.151  1.00  0.00
C
ATOM    193  HA  CYS A  13      -5.382  20.290  29.618  1.00  0.00
H
ATOM    194  CB  CYS A  13      -6.834  21.777  30.224  1.00  0.00
C
ATOM    195  HB1 CYS A  13      -7.474  22.531  29.763  1.00  0.00
H
ATOM    196  HB2 CYS A  13      -6.151  22.319  30.879  1.00  0.00
H
ATOM    197  SG  CYS A  13      -7.869  20.685  31.240  1.00  0.00
S
ATOM    198  C   CYS A  13      -7.005  20.344  28.172  1.00  0.00
C
ATOM    199  O   CYS A  13      -8.022  20.938  27.853  1.00  0.00
O
ATOM    200  N   TYR A  14      -6.677  19.131  27.685  1.00  0.00
N
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ATOM    201  HN  TYR A  14      -7.332  18.848  26.983  1.00  0.00
H
ATOM    202  CA  TYR A  14      -5.578  18.247  28.066  1.00  0.00
C
ATOM    203  HA  TYR A  14      -5.578  18.138  29.142  1.00  0.00
H
ATOM    204  CB  TYR A  14      -5.858  16.873  27.420  1.00  0.00
C
ATOM    205  HB1 TYR A  14      -6.027  16.982  26.354  1.00  0.00
H
ATOM    206  HB2 TYR A  14      -4.983  16.235  27.501  1.00  0.00
H
ATOM    207  CG  TYR A  14      -7.052  16.110  27.953  1.00  0.00
C
ATOM    208  CD1 TYR A  14      -8.031  16.677  28.775  1.00  0.00
C
ATOM    209  HD1 TYR A  14      -7.992  17.701  29.111  1.00  0.00
H
ATOM    210  CE1 TYR A  14      -9.117  15.933  29.210  1.00  0.00
C
ATOM    211  HE1 TYR A  14      -9.865  16.386  29.848  1.00  0.00
H
ATOM    212  CD2 TYR A  14      -7.191  14.771  27.595  1.00  0.00
C
ATOM    213  HD2 TYR A  14      -6.443  14.292  26.972  1.00  0.00
H
ATOM    214  CE2 TYR A  14      -8.286  14.028  28.025  1.00  0.00
C
ATOM    215  HE2 TYR A  14      -8.399  12.993  27.735  1.00  0.00
H
ATOM    216  CZ  TYR A  14      -9.255  14.610  28.828  1.00  0.00
C
ATOM    217  OH  TYR A  14     -10.349  13.879  29.235  1.00  0.00
O
ATOM    218  HH  TYR A  14     -10.337  13.814  30.187  1.00  0.00
H
ATOM    219  C   TYR A  14      -4.240  18.893  27.638  1.00  0.00
C
ATOM    220  O   TYR A  14      -4.239  19.818  26.834  1.00  0.00
O
ATOM    221  N   GLY A  15      -3.114  18.411  28.197  1.00  0.00
N
ATOM    222  HN  GLY A  15      -3.225  17.662  28.856  1.00  0.00
H
ATOM    223  CA  GLY A  15      -1.802  18.995  27.873  1.00  0.00
C
ATOM    224  HA1 GLY A  15      -1.058  18.362  28.351  1.00  0.00
H
ATOM    225  HA2 GLY A  15      -1.762  19.989  28.312  1.00  0.00
H
ATOM    226  C   GLY A  15      -1.559  19.038  26.349  1.00  0.00
C
ATOM    227  O   GLY A  15      -1.610  17.998  25.708  1.00  0.00
O
ATOM    228  N   GLY A  16      -1.397  20.268  25.817  1.00  0.00
N
ATOM    229  HN  GLY A  16      -1.025  20.974  26.418  1.00  0.00
H
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ATOM    230  CA  GLY A  16      -1.870  20.546  24.466  1.00  0.00
C
ATOM    231  HA1 GLY A  16      -2.854  20.089  24.354  1.00  0.00
H
ATOM    232  HA2 GLY A  16      -2.006  21.619  24.389  1.00  0.00
H
ATOM    233  C   GLY A  16      -0.931  20.047  23.363  1.00  0.00
C
ATOM    234  O   GLY A  16       0.262  19.825  23.538  1.00  0.00
O
ATOM    235  N   SER A  17      -1.558  19.868  22.174  1.00  0.00
N
ATOM    236  HN  SER A  17      -2.471  20.261  22.029  1.00  0.00
H
ATOM    237  CA  SER A  17      -0.841  19.162  21.086  1.00  0.00
C
ATOM    238  HA  SER A  17      -0.633  18.163  21.475  1.00  0.00
H
ATOM    239  CB  SER A  17      -1.767  19.058  19.844  1.00  0.00
C
ATOM    240  HB1 SER A  17      -2.773  18.783  20.161  1.00  0.00
H
ATOM    241  HB2 SER A  17      -1.874  20.053  19.417  1.00  0.00
H
ATOM    242  OG  SER A  17      -1.414  18.143  18.801  1.00  0.00
O
ATOM    243  HG  SER A  17      -0.466  17.923  18.857  1.00  0.00
H
ATOM    244  C   SER A  17       0.504  19.861  20.691  1.00  0.00
C
ATOM    245  O   SER A  17       0.516  21.014  20.252  1.00  0.00
O
ATOM    246  N   TYR A  18       1.599  19.060  20.819  1.00  0.00
N
ATOM    247  HN  TYR A  18       1.470  18.144  21.202  1.00  0.00
H
ATOM    248  CA  TYR A  18       2.959  19.485  20.477  1.00  0.00
C
ATOM    249  HA  TYR A  18       3.078  20.522  20.793  1.00  0.00
H
ATOM    250  CB  TYR A  18       3.995  18.519  21.115  1.00  0.00
C
ATOM    251  HB1 TYR A  18       3.560  17.530  21.275  1.00  0.00
H
ATOM    252  HB2 TYR A  18       4.832  18.355  20.434  1.00  0.00
H
ATOM    253  CG  TYR A  18       4.583  19.014  22.415  1.00  0.00
C
ATOM    254  CD1 TYR A  18       3.979  20.028  23.135  1.00  0.00
C
ATOM    255  HD1 TYR A  18       3.015  20.383  22.820  1.00  0.00
H
ATOM    256  CE1 TYR A  18       4.557  20.540  24.279  1.00  0.00
C
ATOM    257  HE1 TYR A  18       4.024  21.302  24.823  1.00  0.00
H
ATOM    258  CD2 TYR A  18       5.759  18.475  22.921  1.00  0.00
C
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ATOM    259  HD2 TYR A  18       6.224  17.646  22.426  1.00  0.00
H
ATOM    260  CE2 TYR A  18       6.340  18.969  24.085  1.00  0.00
C
ATOM    261  HE2 TYR A  18       7.251  18.528  24.458  1.00  0.00
H
ATOM    262  CZ  TYR A  18       5.744  20.017  24.764  1.00  0.00
C
ATOM    263  OH  TYR A  18       6.298  20.502  25.935  1.00  0.00
O
ATOM    264  HH  TYR A  18       6.330  21.460  25.914  1.00  0.00
H
ATOM    265  C   TYR A  18       3.066  19.559  18.970  1.00  0.00
C
ATOM    266  O   TYR A  18       2.205  19.031  18.271  1.00  0.00
O
ATOM    267  N   GLY A  19       4.103  20.303  18.540  1.00  0.00
N
ATOM    268  HN  GLY A  19       4.703  20.583  19.292  1.00  0.00
H
ATOM    269  CA  GLY A  19       4.089  20.956  17.235  1.00  0.00
C
ATOM    270  HA1 GLY A  19       3.161  21.524  17.152  1.00  0.00
H
ATOM    271  HA2 GLY A  19       4.916  21.668  17.254  1.00  0.00
H
ATOM    272  C   GLY A  19       4.238  20.010  16.031  1.00  0.00
C
ATOM    273  O   GLY A  19       3.566  18.982  15.869  1.00  0.00
O
ATOM    274  N   TYR A  20       5.144  20.467  15.131  1.00  0.00
N
ATOM    275  HN  TYR A  20       5.573  21.379  15.231  1.00  0.00
H
ATOM    276  CA  TYR A  20       5.553  19.646  14.005  1.00  0.00
C
ATOM    277  HA  TYR A  20       5.461  18.619  14.355  1.00  0.00
H
ATOM    278  CB  TYR A  20       4.664  19.824  12.747  1.00  0.00
C
ATOM    279  HB1 TYR A  20       5.043  19.167  11.962  1.00  0.00
H
ATOM    280  HB2 TYR A  20       3.671  19.445  12.984  1.00  0.00
H
ATOM    281  CG  TYR A  20       4.515  21.216  12.157  1.00  0.00
C
ATOM    282  CD1 TYR A  20       5.419  21.711  11.216  1.00  0.00
C
ATOM    283  HD1 TYR A  20       6.304  21.141  10.961  1.00  0.00
H
ATOM    284  CE1 TYR A  20       5.198  22.936  10.585  1.00  0.00
C
ATOM    285  HE1 TYR A  20       5.897  23.320   9.852  1.00  0.00
H
ATOM    286  CD2 TYR A  20       3.422  22.010  12.492  1.00  0.00
C
ATOM    287  HD2 TYR A  20       2.735  21.656  13.246  1.00  0.00
H
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ATOM    288  CE2 TYR A  20       3.202  23.238  11.870  1.00  0.00
C
ATOM    289  HE2 TYR A  20       2.348  23.827  12.169  1.00  0.00
H
ATOM    290  CZ  TYR A  20       4.074  23.702  10.886  1.00  0.00
C
ATOM    291  OH  TYR A  20       3.858  24.888  10.180  1.00  0.00
O
ATOM    292  HH  TYR A  20       2.960  25.232  10.131  1.00  0.00
H
ATOM    293  C   TYR A  20       7.025  19.911  13.684  1.00  0.00
C
ATOM    294  O   TYR A  20       7.462  21.062  13.611  1.00  0.00
O
ATOM    295  N   PRO A  21       7.765  18.797  13.376  1.00  0.00
N
ATOM    296  CD  PRO A  21       7.337  17.395  13.467  1.00  0.00
C
ATOM    297  HD1 PRO A  21       7.332  17.086  14.514  1.00  0.00
H
ATOM    298  HD2 PRO A  21       6.349  17.241  13.034  1.00  0.00
H
ATOM    299  CA  PRO A  21       9.049  18.935  12.727  1.00  0.00
C
ATOM    300  HA  PRO A  21       9.680  19.594  13.303  1.00  0.00
H
ATOM    301  CB  PRO A  21       9.592  17.517  12.562  1.00  0.00
C
ATOM    302  HB1 PRO A  21      10.261  17.313  13.388  1.00  0.00
H
ATOM    303  HB2 PRO A  21      10.163  17.355  11.643  1.00  0.00
H
ATOM    304  CG  PRO A  21       8.374  16.605  12.659  1.00  0.00
C
ATOM    305  HG1 PRO A  21       7.999  16.388  11.658  1.00  0.00
H
ATOM    306  HG2 PRO A  21       8.634  15.659  13.126  1.00  0.00
H
ATOM    307  C   PRO A  21       8.747  19.517  11.371  1.00  0.00
C
ATOM    308  O   PRO A  21       7.679  19.219  10.822  1.00  0.00
O
ATOM    309  N   LEU A  22       9.687  20.348  10.898  1.00  0.00
N
ATOM    310  HN  LEU A  22      10.516  20.540  11.431  1.00  0.00
H
ATOM    311  CA  LEU A  22       9.595  20.772   9.519  1.00  0.00
C
ATOM    312  HA  LEU A  22       8.560  21.049   9.313  1.00  0.00
H
ATOM    313  CB  LEU A  22      10.519  21.979   9.249  1.00  0.00
C
ATOM    314  HB1 LEU A  22      10.362  22.337   8.229  1.00  0.00
H
ATOM    315  HB2 LEU A  22      11.562  21.659   9.294  1.00  0.00
H
ATOM    316  CG  LEU A  22      10.315  23.157  10.220  1.00  0.00
C
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ATOM    317  HG  LEU A  22      10.662  22.845  11.204  1.00  0.00
H
ATOM    318  CD1 LEU A  22      11.158  24.369   9.790  1.00  0.00
C
ATOM    319 1HD1 LEU A  22      10.773  24.820   8.873  1.00  0.00
H
ATOM    320 2HD1 LEU A  22      11.147  25.137  10.566  1.00  0.00
H
ATOM    321 3HD1 LEU A  22      12.195  24.091   9.604  1.00  0.00
H
ATOM    322  CD2 LEU A  22       8.841  23.552  10.374  1.00  0.00
C
ATOM    323 1HD2 LEU A  22       8.248  22.743  10.800  1.00  0.00
H
ATOM    324 2HD2 LEU A  22       8.398  23.845   9.421  1.00  0.00
H
ATOM    325 3HD2 LEU A  22       8.744  24.410  11.042  1.00  0.00
H
ATOM    326  C   LEU A  22       9.971  19.592   8.619  1.00  0.00
C
ATOM    327  O   LEU A  22      10.393  18.505   9.022  1.00  0.00
O
ATOM    328  N   GLY A  23       9.846  19.894   7.320  1.00  0.00
N
ATOM    329  HN  GLY A  23       9.454  20.771   7.040  1.00  0.00
H
ATOM    330  CA  GLY A  23      10.483  19.027   6.351  1.00  0.00
C
ATOM    331  HA1 GLY A  23      10.169  17.996   6.523  1.00  0.00
H
ATOM    332  HA2 GLY A  23      10.171  19.344   5.355  1.00  0.00
H
ATOM    333  C   GLY A  23      11.994  19.174   6.505  1.00  0.00
C
ATOM    334  O   GLY A  23      12.508  20.178   7.005  1.00  0.00
O
ATOM    335  N   TYR A  24      12.652  18.100   6.036  1.00  0.00
N
ATOM    336  HN  TYR A  24      12.179  17.313   5.641  1.00  0.00
H
ATOM    337  CA  TYR A  24      14.062  18.180   5.737  1.00  0.00
C
ATOM    338  HA  TYR A  24      14.582  18.371   6.679  1.00  0.00
H
ATOM    339  CB  TYR A  24      14.516  16.827   5.146  1.00  0.00
C
ATOM    340  HB1 TYR A  24      14.467  16.058   5.918  1.00  0.00
H
ATOM    341  HB2 TYR A  24      15.575  16.884   4.890  1.00  0.00
H
ATOM    342  CG  TYR A  24      13.743  16.370   3.925  1.00  0.00
C
ATOM    343  CD1 TYR A  24      14.100  16.816   2.660  1.00  0.00
C
ATOM    344  HD1 TYR A  24      14.912  17.521   2.570  1.00  0.00
H
ATOM    345  CE1 TYR A  24      13.422  16.391   1.524  1.00  0.00
C
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ATOM    346  HE1 TYR A  24      13.727  16.779   0.567  1.00  0.00
H
ATOM    347  CD2 TYR A  24      12.679  15.479   4.028  1.00  0.00
C
ATOM    348  HD2 TYR A  24      12.390  15.096   4.993  1.00  0.00
H
ATOM    349  CE2 TYR A  24      12.003  15.044   2.892  1.00  0.00
C
ATOM    350  HE2 TYR A  24      11.216  14.320   3.005  1.00  0.00
H
ATOM    351  CZ  TYR A  24      12.369  15.489   1.626  1.00  0.00
C
ATOM    352  OH  TYR A  24      11.707  15.023   0.494  1.00  0.00
O
ATOM    353  HH  TYR A  24      12.334  14.864  -0.208  1.00  0.00
H
ATOM    354  C   TYR A  24      14.307  19.341   4.747  1.00  0.00
C
ATOM    355  O   TYR A  24      13.680  19.465   3.698  1.00  0.00
O
ATOM    356  N   SER A  25      15.287  20.171   5.101  1.00  0.00
N
ATOM    357  HN  SER A  25      15.635  20.136   6.041  1.00  0.00
H
ATOM    358  CA  SER A  25      15.952  20.857   4.017  1.00  0.00
C
ATOM    359  HA  SER A  25      15.180  21.328   3.403  1.00  0.00
H
ATOM    360  CB  SER A  25      16.848  21.964   4.591  1.00  0.00
C
ATOM    361  HB1 SER A  25      17.805  21.589   4.952  1.00  0.00
H
ATOM    362  HB2 SER A  25      17.107  22.658   3.791  1.00  0.00
H
ATOM    363  OG  SER A  25      16.178  22.709   5.608  1.00  0.00
O
ATOM    364  HG  SER A  25      16.086  22.196   6.422  1.00  0.00
H
ATOM    365  C   SER A  25      16.662  19.763   3.174  1.00  0.00
C
ATOM    366  O   SER A  25      16.893  18.647   3.628  1.00  0.00
O
ATOM    367  N   VAL A  26      16.902  20.138   1.904  1.00  0.00
N
ATOM    368  HN  VAL A  26      16.691  21.069   1.608  1.00  0.00
H
ATOM    369  CA  VAL A  26      17.587  19.222   1.001  1.00  0.00
C
ATOM    370  HA  VAL A  26      16.957  18.334   0.917  1.00  0.00
H
ATOM    371  CB  VAL A  26      17.775  19.885  -0.378  1.00  0.00
C
ATOM    372  HB  VAL A  26      18.369  20.792  -0.253  1.00  0.00
H
ATOM    373  CG1 VAL A  26      18.528  18.971  -1.358  1.00  0.00
C
ATOM    374 1HG1 VAL A  26      19.523  18.719  -0.989  1.00  0.00
H
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ATOM    375 2HG1 VAL A  26      18.658  19.458  -2.325  1.00  0.00
H
ATOM    376 3HG1 VAL A  26      17.987  18.039  -1.522  1.00  0.00
H
ATOM    377  CG2 VAL A  26      16.418  20.288  -0.977  1.00  0.00
C
ATOM    378 1HG2 VAL A  26      15.888  21.001  -0.345  1.00  0.00
H
ATOM    379 2HG2 VAL A  26      16.542  20.748  -1.958  1.00  0.00
H
ATOM    380 3HG2 VAL A  26      15.773  19.417  -1.096  1.00  0.00
H
ATOM    381  C   VAL A  26      18.922  18.846   1.657  1.00  0.00
C
ATOM    382  O   VAL A  26      19.702  19.707   2.057  1.00  0.00
O
ATOM    383  N   GLY A  27      19.063  17.522   1.840  1.00  0.00
N
ATOM    384  HN  GLY A  27      18.423  16.906   1.378  1.00  0.00
H
ATOM    385  CA  GLY A  27      20.167  16.945   2.598  1.00  0.00
C
ATOM    386  HA1 GLY A  27      20.084  15.862   2.508  1.00  0.00
H
ATOM    387  HA2 GLY A  27      21.098  17.259   2.124  1.00  0.00
H
ATOM    388  C   GLY A  27      20.208  17.315   4.096  1.00  0.00
C
ATOM    389  O   GLY A  27      21.299  17.639   4.582  1.00  0.00
O
ATOM    390  N   CYS A  28      19.024  17.175   4.746  1.00  0.00
N
ATOM    391  HN  CYS A  28      18.182  17.020   4.224  1.00  0.00
H
ATOM    392  CA  CYS A  28      18.837  17.248   6.186  1.00  0.00
C
ATOM    393  HA  CYS A  28      19.810  17.131   6.667  1.00  0.00
H
ATOM    394  CB  CYS A  28      18.188  18.590   6.526  1.00  0.00
C
ATOM    395  HB1 CYS A  28      17.204  18.649   6.067  1.00  0.00
H
ATOM    396  HB2 CYS A  28      18.024  18.701   7.595  1.00  0.00
H
ATOM    397  SG  CYS A  28      19.188  19.976   5.948  1.00  0.00
S
ATOM    398  C   CYS A  28      17.932  16.077   6.596  1.00  0.00
C
ATOM    399  O   CYS A  28      17.458  15.312   5.769  1.00  0.00
O
ATOM    400  N   GLY A  29      17.713  15.952   7.919  1.00  0.00
N
ATOM    401  HN  GLY A  29      17.835  16.707   8.565  1.00  0.00
H
ATOM    402  CA  GLY A  29      17.094  14.716   8.404  1.00  0.00
C
ATOM    403  HA1 GLY A  29      17.313  13.887   7.727  1.00  0.00
H
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ATOM    404  HA2 GLY A  29      17.543  14.488   9.369  1.00  0.00
H
ATOM    405  C   GLY A  29      15.581  14.845   8.593  1.00  0.00
C
ATOM    406  O   GLY A  29      14.999  15.927   8.691  1.00  0.00
O
ATOM    407  N   TYR A  30      14.946  13.654   8.708  1.00  0.00
N
ATOM    408  HN  TYR A  30      15.507  12.868   8.424  1.00  0.00
H
ATOM    409  CA  TYR A  30      13.588  13.666   9.212  1.00  0.00
C
ATOM    410  HA  TYR A  30      13.093  14.392   8.569  1.00  0.00
H
ATOM    411  CB  TYR A  30      12.814  12.352   8.984  1.00  0.00
C
ATOM    412  HB1 TYR A  30      12.773  12.192   7.915  1.00  0.00
H
ATOM    413  HB2 TYR A  30      11.783  12.461   9.319  1.00  0.00
H
ATOM    414  CG  TYR A  30      13.362  11.080   9.568  1.00  0.00
C
ATOM    415  CD1 TYR A  30      12.879  10.614  10.781  1.00  0.00
C
ATOM    416  HD1 TYR A  30      12.184  11.234  11.309  1.00  0.00
H
ATOM    417  CE1 TYR A  30      13.283   9.388  11.284  1.00  0.00
C
ATOM    418  HE1 TYR A  30      12.882   9.007  12.216  1.00  0.00
H
ATOM    419  CD2 TYR A  30      14.299  10.315   8.882  1.00  0.00
C
ATOM    420  HD2 TYR A  30      14.722  10.655   7.948  1.00  0.00
H
ATOM    421  CE2 TYR A  30      14.722   9.096   9.383  1.00  0.00
C
ATOM    422  HE2 TYR A  30      15.456   8.521   8.831  1.00  0.00
H
ATOM    423  CZ  TYR A  30      14.206   8.629  10.580  1.00  0.00
C
ATOM    424  OH  TYR A  30      14.607   7.422  11.098  1.00  0.00
O
ATOM    425  HH  TYR A  30      15.064   7.657  11.906  1.00  0.00
H
ATOM    426  C   TYR A  30      13.574  14.243  10.645  1.00  0.00
C
ATOM    427  O   TYR A  30      13.981  13.638  11.637  1.00  0.00
O
ATOM    428  N   GLY A  31      13.097  15.503  10.693  1.00  0.00
N
ATOM    429  HN  GLY A  31      13.261  15.933   9.802  1.00  0.00
H
ATOM    430  CA  GLY A  31      12.975  16.167  11.987  1.00  0.00
C
ATOM    431  HA1 GLY A  31      13.978  16.239  12.417  1.00  0.00
H
ATOM    432  HA2 GLY A  31      12.581  17.167  11.809  1.00  0.00
H
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ATOM    433  C   GLY A  31      12.027  15.385  12.912  1.00  0.00
C
ATOM    434  O   GLY A  31      11.132  14.661  12.459  1.00  0.00
O
ATOM    435  N   SER A  32      12.210  15.583  14.234  1.00  0.00
N
ATOM    436  HN  SER A  32      12.910  16.257  14.456  1.00  0.00
H
ATOM    437  CA  SER A  32      11.550  14.769  15.258  1.00  0.00
C
ATOM    438  HA  SER A  32      11.396  13.796  14.810  1.00  0.00
H
ATOM    439  CB  SER A  32      12.428  14.590  16.517  1.00  0.00
C
ATOM    440  HB1 SER A  32      13.310  14.006  16.232  1.00  0.00
H
ATOM    441  HB2 SER A  32      12.818  15.549  16.879  1.00  0.00
H
ATOM    442  OG  SER A  32      11.751  13.861  17.550  1.00  0.00
O
ATOM    443  HG  SER A  32      11.379  14.418  18.249  1.00  0.00
H
ATOM    444  C   SER A  32      10.165  15.278  15.652  1.00  0.00
C
ATOM    445  O   SER A  32       9.757  16.429  15.512  1.00  0.00
O
ATOM    446  N   THR A  33       9.419  14.313  16.238  1.00  0.00
N
ATOM    447  HN  THR A  33       9.923  13.496  16.534  1.00  0.00
H
ATOM    448  CA  THR A  33       8.152  14.624  16.901  1.00  0.00
C
ATOM    449  HA  THR A  33       7.658  15.413  16.334  1.00  0.00
H
ATOM    450  CB  THR A  33       7.293  13.330  16.876  1.00  0.00
C
ATOM    451  HB  THR A  33       7.668  12.765  16.024  1.00  0.00
H
ATOM    452  OG1 THR A  33       5.933  13.623  16.571  1.00  0.00
O
ATOM    453  HG1 THR A  33       5.397  12.826  16.635  1.00  0.00
H
ATOM    454  CG2 THR A  33       7.458  12.409  18.101  1.00  0.00
C
ATOM    455 1HG2 THR A  33       8.512  12.297  18.360  1.00  0.00
H
ATOM    456 2HG2 THR A  33       6.935  12.783  18.981  1.00  0.00
H
ATOM    457 3HG2 THR A  33       7.083  11.407  17.896  1.00  0.00
H
ATOM    458  C   THR A  33       8.570  15.176  18.286  1.00  0.00
C
ATOM    459  O   THR A  33       9.764  15.418  18.487  1.00  0.00
O
ATOM    460  N   TYR A  34       7.610  15.310  19.217  1.00  0.00
N
ATOM    461  HN  TYR A  34       6.642  15.095  19.068  1.00  0.00
H
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ATOM    462  CA  TYR A  34       7.952  15.607  20.605  1.00  0.00
C
ATOM    463  HA  TYR A  34       8.840  15.027  20.856  1.00  0.00
H
ATOM    464  CB  TYR A  34       8.201  17.123  20.699  1.00  0.00
C
ATOM    465  HB1 TYR A  34       8.803  17.396  19.835  1.00  0.00
H
ATOM    466  HB2 TYR A  34       7.269  17.667  20.561  1.00  0.00
H
ATOM    467  CG  TYR A  34       8.994  17.707  21.851  1.00  0.00
C
ATOM    468  CD1 TYR A  34       9.790  17.001  22.753  1.00  0.00
C
ATOM    469  HD1 TYR A  34       9.838  15.924  22.727  1.00  0.00
H
ATOM    470  CE1 TYR A  34      10.560  17.683  23.698  1.00  0.00
C
ATOM    471  HE1 TYR A  34      11.154  17.131  24.410  1.00  0.00
H
ATOM    472  CD2 TYR A  34       8.962  19.085  21.963  1.00  0.00
C
ATOM    473  HD2 TYR A  34       8.272  19.616  21.329  1.00  0.00
H
ATOM    474  CE2 TYR A  34       9.757  19.770  22.862  1.00  0.00
C
ATOM    475  HE2 TYR A  34       9.743  20.848  22.864  1.00  0.00
H
ATOM    476  CZ  TYR A  34      10.572  19.075  23.736  1.00  0.00
C
ATOM    477  OH  TYR A  34      11.370  19.771  24.631  1.00  0.00
O
ATOM    478  HH  TYR A  34      10.790  20.106  25.318  1.00  0.00
H
ATOM    479  C   TYR A  34       6.756  15.079  21.415  1.00  0.00
C
ATOM    480  O   TYR A  34       5.791  14.585  20.842  1.00  0.00
O
ATOM    481  N   SER A  35       6.853  15.201  22.741  1.00  0.00
N
ATOM    482  HN  SER A  35       7.656  15.678  23.104  1.00  0.00
H
ATOM    483  CA  SER A  35       6.047  14.526  23.758  1.00  0.00
C
ATOM    484  HA  SER A  35       6.135  13.463  23.532  1.00  0.00
H
ATOM    485  CB  SER A  35       6.764  14.899  25.066  1.00  0.00
C
ATOM    486  HB1 SER A  35       6.471  14.240  25.894  1.00  0.00
H
ATOM    487  HB2 SER A  35       6.449  15.901  25.363  1.00  0.00
H
ATOM    488  OG  SER A  35       8.180  14.956  24.878  1.00  0.00
O
ATOM    489  HG  SER A  35       8.565  14.093  25.074  1.00  0.00
H
ATOM    490  C   SER A  35       4.535  14.934  23.914  1.00  0.00
C
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ATOM    491  O   SER A  35       4.044  15.940  23.400  1.00  0.00
O
ATOM    492  N   PRO A  36       3.823  14.102  24.734  1.00  0.00
N
ATOM    493  CD  PRO A  36       4.156  12.696  24.989  1.00  0.00
C
ATOM    494  HD1 PRO A  36       3.877  12.106  24.117  1.00  0.00
H
ATOM    495  HD2 PRO A  36       5.216  12.547  25.182  1.00  0.00
H
ATOM    496  CA  PRO A  36       2.752  14.586  25.590  1.00  0.00
C
ATOM    497  HA  PRO A  36       1.989  15.102  25.007  1.00  0.00
H
ATOM    498  CB  PRO A  36       2.206  13.336  26.295  1.00  0.00
C
ATOM    499  HB1 PRO A  36       1.887  13.521  27.321  1.00  0.00
H
ATOM    500  HB2 PRO A  36       1.344  12.954  25.747  1.00  0.00
H
ATOM    501  CG  PRO A  36       3.338  12.305  26.224  1.00  0.00
C
ATOM    502  HG1 PRO A  36       2.967  11.279  26.196  1.00  0.00
H
ATOM    503  HG2 PRO A  36       3.955  12.409  27.118  1.00  0.00
H
ATOM    504  C   PRO A  36       3.438  15.565  26.553  1.00  0.00
C
ATOM    505  O   PRO A  36       4.609  15.386  26.906  1.00  0.00
O
ATOM    506  N   VAL A  37       2.692  16.655  26.795  1.00  0.00
N
ATOM    507  HN  VAL A  37       1.744  16.753  26.504  1.00  0.00
H
ATOM    508  CA  VAL A  37       3.213  17.742  27.589  1.00  0.00
C
ATOM    509  HA  VAL A  37       4.217  17.493  27.942  1.00  0.00
H
ATOM    510  CB  VAL A  37       3.219  19.032  26.774  1.00  0.00
C
ATOM    511  HB  VAL A  37       3.860  18.784  25.938  1.00  0.00
H
ATOM    512  CG1 VAL A  37       1.842  19.422  26.214  1.00  0.00
C
ATOM    513 1HG1 VAL A  37       1.132  19.618  27.017  1.00  0.00
H
ATOM    514 2HG1 VAL A  37       1.424  18.646  25.574  1.00  0.00
H
ATOM    515 3HG1 VAL A  37       1.907  20.333  25.619  1.00  0.00
H
ATOM    516  CG2 VAL A  37       3.838  20.197  27.564  1.00  0.00
C
ATOM    517 1HG2 VAL A  37       3.157  20.575  28.326  1.00  0.00
H
ATOM    518 2HG2 VAL A  37       4.775  19.910  28.041  1.00  0.00
H
ATOM    519 3HG2 VAL A  37       4.055  21.049  26.934  1.00  0.00
H
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ATOM    520  C   VAL A  37       2.301  17.889  28.769  1.00  0.00
C
ATOM    521  O   VAL A  37       1.166  17.417  28.774  1.00  0.00
O
ATOM    522  N   GLY A  38       2.890  18.539  29.799  1.00  0.00
N
ATOM    523  HN  GLY A  38       3.832  18.831  29.603  1.00  0.00
H
ATOM    524  CA  GLY A  38       2.204  19.001  31.003  1.00  0.00
C
ATOM    525  HA1 GLY A  38       2.636  19.968  31.259  1.00  0.00
H
ATOM    526  HA2 GLY A  38       2.434  18.288  31.797  1.00  0.00
H
ATOM    527  C   GLY A  38       0.687  19.158  30.802  1.00  0.00
C
ATOM    528  O   GLY A  38       0.217  19.890  29.927  1.00  0.00
O
ATOM    529  N   TYR A  39       0.028  18.342  31.624  1.00  0.00
N
ATOM    530  HN  TYR A  39       0.540  17.865  32.339  1.00  0.00
H
ATOM    531  CA  TYR A  39      -1.394  18.137  31.678  1.00  0.00
C
ATOM    532  HA  TYR A  39      -1.883  18.776  30.941  1.00  0.00
H
ATOM    533  CB  TYR A  39      -1.677  16.641  31.422  1.00  0.00
C
ATOM    534  HB1 TYR A  39      -1.621  16.464  30.349  1.00  0.00
H
ATOM    535  HB2 TYR A  39      -2.696  16.405  31.735  1.00  0.00
H
ATOM    536  CG  TYR A  39      -0.725  15.678  32.104  1.00  0.00
C
ATOM    537  CD1 TYR A  39      -0.900  15.347  33.443  1.00  0.00
C
ATOM    538  HD1 TYR A  39      -1.720  15.799  33.986  1.00  0.00
H
ATOM    539  CE1 TYR A  39      -0.043  14.458  34.079  1.00  0.00
C
ATOM    540  HE1 TYR A  39      -0.220  14.259  35.125  1.00  0.00
H
ATOM    541  CD2 TYR A  39       0.333  15.090  31.411  1.00  0.00
C
ATOM    542  HD2 TYR A  39       0.483  15.342  30.372  1.00  0.00
H
ATOM    543  CE2 TYR A  39       1.185  14.185  32.038  1.00  0.00
C
ATOM    544  HE2 TYR A  39       1.992  13.698  31.509  1.00  0.00
H
ATOM    545  CZ  TYR A  39       1.003  13.863  33.376  1.00  0.00
C
ATOM    546  OH  TYR A  39       1.852  12.942  33.983  1.00  0.00
O
ATOM    547  HH  TYR A  39       1.486  12.592  34.796  1.00  0.00
H
ATOM    548  C   TYR A  39      -1.784  18.586  33.077  1.00  0.00
C
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ATOM    549  O   TYR A  39      -1.097  18.287  34.047  1.00  0.00
O
ATOM    550  N   GLY A  40      -2.873  19.360  33.180  1.00  0.00
N
ATOM    551  HN  GLY A  40      -3.306  19.571  32.304  1.00  0.00
H
ATOM    552  CA  GLY A  40      -3.310  19.819  34.499  1.00  0.00
C
ATOM    553  HA1 GLY A  40      -2.459  20.274  35.009  1.00  0.00
H
ATOM    554  HA2 GLY A  40      -4.078  20.573  34.331  1.00  0.00
H
ATOM    555  C   GLY A  40      -3.889  18.663  35.335  1.00  0.00
C
ATOM    556  O   GLY A  40      -3.989  17.528  34.883  1.00  0.00
O
ATOM    557  N   PHE A  41      -4.321  19.011  36.574  1.00  0.00
N
ATOM    558  HN  PHE A  41      -4.237  19.928  36.971  1.00  0.00
H
ATOM    559  CA  PHE A  41      -5.210  18.060  37.252  1.00  0.00
C
ATOM    560  HA  PHE A  41      -4.715  17.086  37.243  1.00  0.00
H
ATOM    561  CB  PHE A  41      -5.507  18.497  38.703  1.00  0.00
C
ATOM    562  HB1 PHE A  41      -4.576  18.663  39.248  1.00  0.00
H
ATOM    563  HB2 PHE A  41      -6.049  19.444  38.721  1.00  0.00
H
ATOM    564  CG  PHE A  41      -6.315  17.449  39.431  1.00  0.00
C
ATOM    565  CD1 PHE A  41      -5.690  16.516  40.254  1.00  0.00
C
ATOM    566  HD1 PHE A  41      -4.625  16.562  40.434  1.00  0.00
H
ATOM    567  CD2 PHE A  41      -7.697  17.377  39.267  1.00  0.00
C
ATOM    568  HD2 PHE A  41      -8.213  18.098  38.644  1.00  0.00
H
ATOM    569  CE1 PHE A  41      -6.421  15.510  40.870  1.00  0.00
C
ATOM    570  HE1 PHE A  41      -5.927  14.788  41.507  1.00  0.00
H
ATOM    571  CE2 PHE A  41      -8.430  16.369  39.879  1.00  0.00
C
ATOM    572  HE2 PHE A  41      -9.499  16.312  39.728  1.00  0.00
H
ATOM    573  CZ  PHE A  41      -7.791  15.433  40.676  1.00  0.00
C
ATOM    574  HZ  PHE A  41      -8.354  14.649  41.164  1.00  0.00
H
ATOM    575  C   PHE A  41      -6.518  17.988  36.438  1.00  0.00
C
ATOM    576  O   PHE A  41      -7.002  19.005  35.947  1.00  0.00
O
ATOM    577  N   GLY A  42      -7.041  16.759  36.281  1.00  0.00
N
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ATOM    578  HN  GLY A  42      -6.485  15.973  36.550  1.00  0.00
H
ATOM    579  CA  GLY A  42      -8.335  16.576  35.625  1.00  0.00
C
ATOM    580  HA1 GLY A  42      -8.999  17.401  35.891  1.00  0.00
H
ATOM    581  HA2 GLY A  42      -8.759  15.650  36.015  1.00  0.00
H
ATOM    582  C   GLY A  42      -8.194  16.470  34.100  1.00  0.00
C
ATOM    583  O   GLY A  42      -8.871  15.677  33.456  1.00  0.00
O
ATOM    584  N   CYS A  43      -7.284  17.320  33.603  1.00  0.00
N
ATOM    585  HN  CYS A  43      -6.904  17.997  34.229  1.00  0.00
H
ATOM    586  CA  CYS A  43      -6.787  17.271  32.256  1.00  0.00
C
ATOM    587  HA  CYS A  43      -7.642  17.286  31.581  1.00  0.00
H
ATOM    588  CB  CYS A  43      -5.889  18.502  32.107  1.00  0.00
C
ATOM    589  HB1 CYS A  43      -5.553  18.633  31.088  1.00  0.00
H
ATOM    590  HB2 CYS A  43      -4.995  18.310  32.688  1.00  0.00
H
ATOM    591  SG  CYS A  43      -6.601  20.058  32.711  1.00  0.00
S
ATOM    592  C   CYS A  43      -5.949  15.990  32.103  1.00  0.00
C
ATOM    593  O   CYS A  43      -5.128  15.662  32.947  1.00  0.00
O
ATOM    594  N   GLY A  44      -6.184  15.283  30.993  1.00  0.00
N
ATOM    595  HN  GLY A  44      -6.958  15.520  30.411  1.00  0.00
H
ATOM    596  CA  GLY A  44      -5.241  14.245  30.602  1.00  0.00
C
ATOM    597  HA1 GLY A  44      -5.787  13.523  29.994  1.00  0.00
H
ATOM    598  HA2 GLY A  44      -4.846  13.737  31.486  1.00  0.00
H
ATOM    599  C   GLY A  44      -4.113  14.881  29.786  1.00  0.00
C
ATOM    600  O   GLY A  44      -3.981  16.101  29.686  1.00  0.00
O
ATOM    601  N   ALA A  45      -3.323  13.981  29.165  1.00  0.00
N
ATOM    602  HN  ALA A  45      -3.556  13.006  29.242  1.00  0.00
H
ATOM    603  CA  ALA A  45      -2.352  14.459  28.171  1.00  0.00
C
ATOM    604  HA  ALA A  45      -2.049  15.479  28.411  1.00  0.00
H
ATOM    605  CB  ALA A  45      -1.131  13.531  28.175  1.00  0.00
C
ATOM    606  HB1 ALA A  45      -0.370  13.897  27.488  1.00  0.00
H
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ATOM    607  HB2 ALA A  45      -1.403  12.520  27.866  1.00  0.00
H
ATOM    608  HB3 ALA A  45      -0.691  13.469  29.170  1.00  0.00
H
ATOM    609  C   ALA A  45      -3.050  14.442  26.803  1.00  0.00
C
ATOM    610  O   ALA A  45      -4.047  13.722  26.673  1.00  0.00
O
ATOM    611  N   PHE A  46      -2.480  15.251  25.850  1.00  0.00
N
ATOM    612  HN  PHE A  46      -1.867  15.977  26.166  1.00  0.00
H
ATOM    613  CA  PHE A  46      -2.704  15.002  24.416  1.00  0.00
C
ATOM    614  HA  PHE A  46      -3.731  14.648  24.316  1.00  0.00
H
ATOM    615  CB  PHE A  46      -2.492  16.220  23.468  1.00  0.00
C
ATOM    616  HB1 PHE A  46      -1.624  16.804  23.764  1.00  0.00
H
ATOM    617  HB2 PHE A  46      -2.256  15.869  22.461  1.00  0.00
H
ATOM    618  CG  PHE A  46      -3.699  17.108  23.310  1.00  0.00
C
ATOM    619  CD1 PHE A  46      -4.311  17.683  24.414  1.00  0.00
C
ATOM    620  HD1 PHE A  46      -3.912  17.487  25.397  1.00  0.00
H
ATOM    621  CD2 PHE A  46      -4.220  17.372  22.044  1.00  0.00
C
ATOM    622  HD2 PHE A  46      -3.765  16.931  21.169  1.00  0.00
H
ATOM    623  CE1 PHE A  46      -5.418  18.507  24.262  1.00  0.00
C
ATOM    624  HE1 PHE A  46      -5.879  18.949  25.132  1.00  0.00
H
ATOM    625  CE2 PHE A  46      -5.336  18.188  21.891  1.00  0.00
C
ATOM    626  HE2 PHE A  46      -5.764  18.378  20.915  1.00  0.00
H
ATOM    627  CZ  PHE A  46      -5.937  18.755  23.004  1.00  0.00
C
ATOM    628  HZ  PHE A  46      -6.825  19.368  22.903  1.00  0.00
H
ATOM    629  C   PHE A  46      -1.788  13.846  23.970  1.00  0.00
C
ATOM    630  O   PHE A  46      -2.160  12.683  24.077  1.00  0.00
O
ATOM    631  N   GLY A  47      -0.616  14.197  23.394  1.00  0.00
N
ATOM    632  HN  GLY A  47      -0.319  15.152  23.464  1.00  0.00
H
ATOM    633  CA  GLY A  47       0.157  13.175  22.705  1.00  0.00
C
ATOM    634  HA1 GLY A  47      -0.518  12.480  22.202  1.00  0.00
H
ATOM    635  HA2 GLY A  47       0.755  12.645  23.447  1.00  0.00
H
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ATOM    636  C   GLY A  47       1.051  13.857  21.687  1.00  0.00
C
ATOM    637  O   GLY A  47       1.175  15.085  21.675  1.00  0.00
O
ATOM    638  N   TYR A  48       1.711  13.001  20.881  1.00  0.00
N
ATOM    639  HN  TYR A  48       1.547  12.019  20.863  1.00  0.00
H
ATOM    640  CA  TYR A  48       2.732  13.479  19.954  1.00  0.00
C
ATOM    641  HA  TYR A  48       3.108  14.445  20.303  1.00  0.00
H
ATOM    642  CB  TYR A  48       3.891  12.461  19.914  1.00  0.00
C
ATOM    643  HB1 TYR A  48       4.588  12.676  19.106  1.00  0.00
H
ATOM    644  HB2 TYR A  48       4.470  12.570  20.831  1.00  0.00
H
ATOM    645  CG  TYR A  48       3.442  11.023  19.825  1.00  0.00
C
ATOM    646  CD1 TYR A  48       2.804  10.528  18.690  1.00  0.00
C
ATOM    647  HD1 TYR A  48       2.604  11.163  17.846  1.00  0.00
H
ATOM    648  CE1 TYR A  48       2.382   9.210  18.624  1.00  0.00
C
ATOM    649  HE1 TYR A  48       1.899   8.829  17.734  1.00  0.00
H
ATOM    650  CD2 TYR A  48       3.655  10.166  20.901  1.00  0.00
C
ATOM    651  HD2 TYR A  48       4.142  10.509  21.804  1.00  0.00
H
ATOM    652  CE2 TYR A  48       3.225   8.848  20.844  1.00  0.00
C
ATOM    653  HE2 TYR A  48       3.387   8.194  21.684  1.00  0.00
H
ATOM    654  CZ  TYR A  48       2.584   8.375  19.711  1.00  0.00
C
ATOM    655  OH  TYR A  48       2.147   7.071  19.668  1.00  0.00
O
ATOM    656  HH  TYR A  48       2.640   6.610  18.990  1.00  0.00
H
ATOM    657  C   TYR A  48       2.086  13.683  18.579  1.00  0.00
C
ATOM    658  O   TYR A  48       0.922  13.310  18.373  1.00  0.00
O
ATOM    659  N   ARG A  49       2.937  14.273  17.695  1.00  0.00
N
ATOM    660  HN  ARG A  49       3.885  14.463  17.965  1.00  0.00
H
ATOM    661  CA  ARG A  49       2.580  14.476  16.302  1.00  0.00
C
ATOM    662  HA  ARG A  49       1.524  14.224  16.195  1.00  0.00
H
ATOM    663  CB  ARG A  49       2.784  15.976  15.951  1.00  0.00
C
ATOM    664  HB1 ARG A  49       2.668  16.573  16.857  1.00  0.00
H
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ATOM    665  HB2 ARG A  49       3.816  16.155  15.646  1.00  0.00
H
ATOM    666  CG  ARG A  49       1.809  16.543  14.892  1.00  0.00
C
ATOM    667  HG1 ARG A  49       2.304  17.351  14.351  1.00  0.00
H
ATOM    668  HG2 ARG A  49       1.621  15.779  14.136  1.00  0.00
H
ATOM    669  CD  ARG A  49       0.465  17.083  15.450  1.00  0.00
C
ATOM    670  HD1 ARG A  49      -0.058  16.330  16.041  1.00  0.00
H
ATOM    671  HD2 ARG A  49      -0.184  17.371  14.624  1.00  0.00
H
ATOM    672  NE  ARG A  49       0.663  18.269  16.298  1.00  0.00
N1+
ATOM    673  HE  ARG A  49       1.598  18.319  16.700  1.00  0.00
H
ATOM    674  CZ  ARG A  49      -0.269  19.258  16.543  1.00  0.00
C
ATOM    675  NH1 ARG A  49       0.030  20.339  17.272  1.00  0.00
N
ATOM    676 1HH1 ARG A  49       0.921  20.355  17.746  1.00  0.00
H
ATOM    677 2HH1 ARG A  49      -0.602  21.113  17.377  1.00  0.00
H
ATOM    678  NH2 ARG A  49      -1.504  19.128  16.028  1.00  0.00
N
ATOM    679 1HH2 ARG A  49      -1.744  18.317  15.485  1.00  0.00
H
ATOM    680 2HH2 ARG A  49      -2.257  19.782  16.144  1.00  0.00
H
ATOM    681  C   ARG A  49       3.398  13.418  15.512  1.00  0.00
C
ATOM    682  O   ARG A  49       3.350  12.233  15.862  1.00  0.00
O
ATOM    683  N   ARG A  50       4.098  13.851  14.453  1.00  0.00
N
ATOM    684  HN  ARG A  50       4.345  14.815  14.361  1.00  0.00
H
ATOM    685  CA  ARG A  50       4.854  12.938  13.628  1.00  0.00
C
ATOM    686  HA  ARG A  50       4.707  11.904  13.952  1.00  0.00
H
ATOM    687  CB  ARG A  50       4.370  13.148  12.183  1.00  0.00
C
ATOM    688  HB1 ARG A  50       4.712  14.117  11.814  1.00  0.00
H
ATOM    689  HB2 ARG A  50       4.829  12.385  11.552  1.00  0.00
H
ATOM    690  CG  ARG A  50       2.844  13.092  12.044  1.00  0.00
C
ATOM    691  HG1 ARG A  50       2.451  12.351  12.744  1.00  0.00
H
ATOM    692  HG2 ARG A  50       2.402  14.047  12.331  1.00  0.00
H
ATOM    693  CD  ARG A  50       2.386  12.689  10.639  1.00  0.00
C
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ATOM    694  HD1 ARG A  50       1.296  12.721  10.599  1.00  0.00
H
ATOM    695  HD2 ARG A  50       2.707  11.668  10.434  1.00  0.00
H
ATOM    696  NE  ARG A  50       2.978  13.530   9.587  1.00  0.00
N1+
ATOM    697  HE  ARG A  50       3.918  13.867   9.786  1.00  0.00
H
ATOM    698  CZ  ARG A  50       2.351  13.763   8.399  1.00  0.00
C
ATOM    699  NH1 ARG A  50       3.056  14.414   7.462  1.00  0.00
N
ATOM    700 1HH1 ARG A  50       4.036  14.660   7.586  1.00  0.00
H
ATOM    701 2HH1 ARG A  50       2.632  14.686   6.592  1.00  0.00
H
ATOM    702  NH2 ARG A  50       1.089  13.347   8.212  1.00  0.00
N
ATOM    703 1HH2 ARG A  50       0.633  12.960   9.018  1.00  0.00
H
ATOM    704 2HH2 ARG A  50       0.581  13.365   7.341  1.00  0.00
H
ATOM    705  C   ARG A  50       6.328  13.320  13.761  1.00  0.00
C
ATOM    706  O   ARG A  50       6.665  14.437  14.142  1.00  0.00
O
ATOM    707  N   TYR A  51       7.171  12.377  13.258  1.00  0.00
N
ATOM    708  HN  TYR A  51       6.875  11.454  12.980  1.00  0.00
H
ATOM    709  CA  TYR A  51       8.454  12.759  12.661  1.00  0.00
C
ATOM    710  HA  TYR A  51       8.913  13.536  13.273  1.00  0.00
H
ATOM    711  CB  TYR A  51       9.371  11.512  12.640  1.00  0.00
C
ATOM    712  HB1 TYR A  51       8.810  10.633  12.319  1.00  0.00
H
ATOM    713  HB2 TYR A  51      10.164  11.641  11.905  1.00  0.00
H
ATOM    714  CG  TYR A  51      10.008  11.233  13.990  1.00  0.00
C
ATOM    715  CD1 TYR A  51       9.251  10.875  15.108  1.00  0.00
C
ATOM    716  HD1 TYR A  51       8.184  10.716  15.045  1.00  0.00
H
ATOM    717  CE1 TYR A  51       9.848  10.708  16.354  1.00  0.00
C
ATOM    718  HE1 TYR A  51       9.220  10.450  17.191  1.00  0.00
H
ATOM    719  CD2 TYR A  51      11.386  11.353  14.156  1.00  0.00
C
ATOM    720  HD2 TYR A  51      12.037  11.632  13.340  1.00  0.00
H
ATOM    721  CE2 TYR A  51      11.986  11.171  15.392  1.00  0.00
C
ATOM    722  HE2 TYR A  51      13.060  11.268  15.520  1.00  0.00
H
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ATOM    723  CZ  TYR A  51      11.219  10.875  16.501  1.00  0.00
C
ATOM    724  OH  TYR A  51      11.883  10.766  17.716  1.00  0.00
O
ATOM    725  HH  TYR A  51      11.313  10.703  18.495  1.00  0.00
H
ATOM    726  C   TYR A  51       8.094  13.362  11.267  1.00  0.00
C
ATOM    727  O   TYR A  51       6.909  13.660  11.063  1.00  0.00
O
ATOM    728  N   SER A  52       9.126  13.552  10.408  1.00  0.00
N
ATOM    729  HN  SER A  52      10.049  13.193  10.590  1.00  0.00
H
ATOM    730  CA  SER A  52       8.979  14.042   9.035  1.00  0.00
C
ATOM    731  HA  SER A  52       8.017  14.550   8.940  1.00  0.00
H
ATOM    732  CB  SER A  52      10.175  15.005   8.770  1.00  0.00
C
ATOM    733  HB1 SER A  52      10.024  15.914   9.361  1.00  0.00
H
ATOM    734  HB2 SER A  52      11.047  14.511   9.201  1.00  0.00
H
ATOM    735  OG  SER A  52      10.441  15.418   7.420  1.00  0.00
O
ATOM    736  HG  SER A  52      11.374  15.659   7.303  1.00  0.00
H
ATOM    737  C   SER A  52       8.998  12.782   8.144  1.00  0.00
C
ATOM    738  O   SER A  52       9.191  11.653   8.618  1.00  0.00
O
ATOM    739  N   PRO A  53       8.771  12.977   6.814  1.00  0.00
N
ATOM    740  CD  PRO A  53       7.982  14.066   6.239  1.00  0.00
C
ATOM    741  HD1 PRO A  53       8.558  14.990   6.211  1.00  0.00
H
ATOM    742  HD2 PRO A  53       7.071  14.242   6.806  1.00  0.00
H
ATOM    743  CA  PRO A  53       9.331  12.082   5.805  1.00  0.00
C
ATOM    744  HA  PRO A  53       8.950  11.071   5.955  1.00  0.00
H
ATOM    745  CB  PRO A  53       8.845  12.659   4.464  1.00  0.00
C
ATOM    746  HB1 PRO A  53       8.573  11.879   3.750  1.00  0.00
H
ATOM    747  HB2 PRO A  53       9.624  13.263   4.003  1.00  0.00
H
ATOM    748  CG  PRO A  53       7.673  13.576   4.822  1.00  0.00
C
ATOM    749  HG1 PRO A  53       7.529  14.394   4.115  1.00  0.00
H
ATOM    750  HG2 PRO A  53       6.760  12.981   4.834  1.00  0.00
H
ATOM    751  C   PRO A  53      10.870  12.085   5.886  1.00  0.00
C
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ATOM    752  O   PRO A  53      11.478  13.102   6.222  1.00  0.00
O
ATOM    753  N   PHE A  54      11.444  10.909   5.552  1.00  0.00
N
ATOM    754  HN  PHE A  54      10.892  10.147   5.216  1.00  0.00
H
ATOM    755  CA  PHE A  54      12.874  10.906   5.274  1.00  0.00
C
ATOM    756  HA  PHE A  54      13.361  11.523   6.028  1.00  0.00
H
ATOM    757  CB  PHE A  54      13.463   9.479   5.375  1.00  0.00
C
ATOM    758  HB1 PHE A  54      14.457   9.568   5.811  1.00  0.00
H
ATOM    759  HB2 PHE A  54      12.901   8.895   6.102  1.00  0.00
H
ATOM    760  CG  PHE A  54      13.591   8.661   4.107  1.00  0.00
C
ATOM    761  CD1 PHE A  54      14.768   7.949   3.877  1.00  0.00
C
ATOM    762  HD1 PHE A  54      15.592   8.019   4.580  1.00  0.00
H
ATOM    763  CD2 PHE A  54      12.561   8.567   3.171  1.00  0.00
C
ATOM    764  HD2 PHE A  54      11.648   9.126   3.312  1.00  0.00
H
ATOM    765  CE1 PHE A  54      14.908   7.144   2.753  1.00  0.00
C
ATOM    766  HE1 PHE A  54      15.823   6.588   2.593  1.00  0.00
H
ATOM    767  CE2 PHE A  54      12.698   7.762   2.044  1.00  0.00
C
ATOM    768  HE2 PHE A  54      11.897   7.674   1.317  1.00  0.00
H
ATOM    769  CZ  PHE A  54      13.871   7.051   1.838  1.00  0.00
C
ATOM    770  HZ  PHE A  54      13.973   6.420   0.965  1.00  0.00
H
ATOM    771  C   PHE A  54      13.045  11.613   3.931  1.00  0.00
C
ATOM    772  O   PHE A  54      12.073  11.758   3.190  1.00  0.00
O
ATOM    773  N   ALA A  55      14.295  12.014   3.703  1.00  0.00
N
ATOM    774  HN  ALA A  55      14.918  11.681   4.408  1.00  0.00
H
ATOM    775  CA  ALA A  55      14.758  12.869   2.623  1.00  0.00
C
ATOM    776  HA  ALA A  55      13.908  13.402   2.216  1.00  0.00
H
ATOM    777  CB  ALA A  55      15.821  13.815   3.216  1.00  0.00
C
ATOM    778  HB1 ALA A  55      16.737  13.287   3.487  1.00  0.00
H
ATOM    779  HB2 ALA A  55      15.472  14.310   4.123  1.00  0.00
H
ATOM    780  HB3 ALA A  55      16.100  14.581   2.494  1.00  0.00
H
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ATOM    781  C   ALA A  55      15.383  12.048   1.493  1.00  0.00
C
ATOM    782  O   ALA A  55      15.691  12.543   0.412  1.00  0.00
O
ATOM    783  N   LEU A  56      15.595  10.767   1.833  1.00  0.00
N
ATOM    784  HN  LEU A  56      15.056  10.415   2.596  1.00  0.00
H
ATOM    785  CA  LEU A  56      16.723   9.939   1.414  1.00  0.00
C
ATOM    786  HA  LEU A  56      16.296   8.959   1.220  1.00  0.00
H
ATOM    787  CB  LEU A  56      17.514  10.394   0.163  1.00  0.00
C
ATOM    788  HB1 LEU A  56      17.848  11.423   0.298  1.00  0.00
H
ATOM    789  HB2 LEU A  56      18.432   9.809   0.078  1.00  0.00
H
ATOM    790  CG  LEU A  56      16.731  10.242  -1.156  1.00  0.00
C
ATOM    791  HG  LEU A  56      15.709  10.589  -0.996  1.00  0.00
H
ATOM    792  CD1 LEU A  56      17.340  11.124  -2.253  1.00  0.00
C
ATOM    793 1HD1 LEU A  56      17.354  12.172  -1.954  1.00  0.00
H
ATOM    794 2HD1 LEU A  56      18.363  10.823  -2.485  1.00  0.00
H
ATOM    795 3HD1 LEU A  56      16.762  11.057  -3.174  1.00  0.00
H
ATOM    796  CD2 LEU A  56      16.664   8.778  -1.620  1.00  0.00
C
ATOM    797 1HD2 LEU A  56      16.160   8.141  -0.893  1.00  0.00
H
ATOM    798 2HD2 LEU A  56      17.656   8.361  -1.800  1.00  0.00
H
ATOM    799 3HD2 LEU A  56      16.105   8.695  -2.553  1.00  0.00
H
ATOM    800  C   LEU A  56      17.656   9.786   2.626  1.00  0.00
C
ATOM    801  O   LEU A  56      18.488   8.883   2.646  1.00  0.00
O
ATOM    802  N   TYR A  57      17.449  10.678   3.623  1.00  0.00
N
ATOM    803  HN  TYR A  57      16.724  11.365   3.614  1.00  0.00
H
ATOM    804  CA  TYR A  57      18.082  10.521   4.919  1.00  0.00
C
ATOM    805  HA  TYR A  57      18.519   9.521   4.974  1.00  0.00
H
ATOM    806  CB  TYR A  57      19.167  11.590   5.114  1.00  0.00
C
ATOM    807  HB1 TYR A  57      19.741  11.712   4.196  1.00  0.00
H
ATOM    808  HB2 TYR A  57      18.723  12.560   5.352  1.00  0.00
H
ATOM    809  CG  TYR A  57      20.092  11.153   6.213  1.00  0.00
C
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ATOM    810  CD1 TYR A  57      21.245  10.431   5.930  1.00  0.00
C
ATOM    811  HD1 TYR A  57      21.533  10.200   4.913  1.00  0.00
H
ATOM    812  CE1 TYR A  57      22.046   9.955   6.960  1.00  0.00
C
ATOM    813  HE1 TYR A  57      22.928   9.398   6.700  1.00  0.00
H
ATOM    814  CD2 TYR A  57      19.771  11.417   7.535  1.00  0.00
C
ATOM    815  HD2 TYR A  57      18.868  11.972   7.750  1.00  0.00
H
ATOM    816  CE2 TYR A  57      20.558  10.934   8.568  1.00  0.00
C
ATOM    817  HE2 TYR A  57      20.250  11.111   9.588  1.00  0.00
H
ATOM    818  CZ  TYR A  57      21.695  10.193   8.283  1.00  0.00
C
ATOM    819  OH  TYR A  57      22.464   9.698   9.322  1.00  0.00
O
ATOM    820  HH  TYR A  57      22.837   8.846   9.115  1.00  0.00
H
ATOM    821  C   TYR A  57      16.967  10.632   5.975  1.00  0.00
C
ATOM    822 1OCT TYR A  57      16.937   9.842   6.915  1.00  0.00
O
ATOM    823 2OCT TYR A  57      16.109  11.498   5.829  1.00  0.00
O1-
TER     824      TYR A  57
CONECT   46  397
CONECT  197  591
CONECT  397   46
CONECT  591  197
END
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