SUPPLEMENTAL INFORMATION

Construction of plasmids used in the study

PEC2003: (GST) The plasmid PEC610 (*UAF30*/pCBGST1) in which yeast *UAF30* is cloned as *Nco*I/*Kpn*I fragment into pCBGST1 was digested with *Nco*I. The *Nco*I site was filled-in with Klenow and dNTP mix and the linear plasmid was ligated. This results in the generation of a stop codon 6 base pairs (two amino acid residues) downstream of the in-frame ATG codon present in the original *Nco*I site.

PEC2007: (GST-proRTD1a) The proRTD1a sequence was amplified by PCR using oligonucleotides #512 and #513 and using plasmid CIC/pLNCX (from Dr. Jun Yuan, this has a mutation of residue $T^{38}A$ in the prosegment) as template. The PCR product was digested with *Nco*I and *Acc*65I and ligated into plasmid PEC610 previously digested with the same restriction enzymes.

PEC2038: (DBD-proRTD1a) The proRTD1a sequence was amplified by PCR using oligonucleotides #506 and #512 using plasmid CIC/pLNCX as template. The PCR product was digested with *Nco*I and *Bam*HI and ligated into plasmid pGBKT7 (Clontech) also digested with *Nco*I and *Bam*HI.

PEC2069: (AD-SDF2L1) This plasmid expresses the rhesus SDF2L1 as a fusion to the GAL4 activation domain produced as a result of the recombination between the amplified cDNA and linearized pGADT7-Rec plasmid. This was isolated from a yeast colony containing a putative DBD-proRTD1a interactor.

PEC2073: (DBD-proseg) The prosegment of proRTD1a was amplified by PCR using oligonucleotides #552 and #553 using PEC2038 as template. The PCR product was digested *Eco*RI and *Bam*HI and ligated into pGBKT7 vector also digested with the same enzymes.

PEC2075: (GST-ΔN_{28}SDF2L1) Oligonucleotides #556 and #557 were used to amplify ΔN_{28}SDF2L1 by PCR using PEC2069 DNA as template. The PCR product was digested with *Bsp*HI and *Acc*65I and ligated into PEC2007 DNA digested with *Nco*I and *Acc*65I.

PEC2082: (DBD-proHNP3) The plasmid PEC2080 (proHNP3/pCR2.1) was first digested with *Acc*65I and the site was then filled-in using Klenow and dNTPs. The plasmid was then digested with *Eco*RI and the proHNP3 containing fragment was ligated into pGBKT7 vector digested with *Eco*RI and *Sma*I.

PEC2084: (AD-ΔN_{28}SDF2L1) Plasmid PEC2069 was used as a template to amplify ΔN_{28}SDF2L1 using oligonucleotides #536 and #556. The PCR product was digested with *Eco*RI and *Xho*I and ligated into plasmid pGADT7-Rec digested with the same enzymes.

PEC2105: (AD-SDF2L1ΔMIR1) Oligonucleotides #557 and #580 were used to amplify SDF2L1-ΔMIR1 (ΔN_{87}-SDF2L1) using PEC2069 as template by PCR. The PCR product

was digested with *Eco*RI and *Xho*I and ligated into pGADT7-Rec digested with the same enzymes.

PEC2106: (AD-SDF2L1ΔMIR3) Oligonucleotides #535 and #581 were used to amplify SDF2L1-ΔMIR3 (SDF2L1-ΔC_{151}) using PEC2069 as template by PCR. The PCR product was digested with *Eco*RI and *Xho*I and ligated into pGADT7-Rec vector also digested with the same enzymes.

PEC2113: (DBD-proHBD1) The proHBD1 fragment was amplified by PCR using oligonucleotides #586 and #587 using PEC2083 (PreproHBD1/pCR2.1 Topo) as template. The PCR product was then cloned into pGBKT7 as *Eco*RI and *Bam*HI fragment.

PEC2114: (DBD-proHD5) The proHD5 fragment was amplified by PCR using oligonucleotides #584 and #585 and plasmid PEC2109 (full length HD5 cDNA, item # IHS1380-97652042 from Open Biosystems) as template and cloned as *Eco*RI and *Bam*HI fragment into pGBKT7.

PEC2121: (GST-proHNP3) The proHNP3 fragment was amplified using PEC2082 as template and oligonucleotides #597 and #570 by PCR. The PCR product was digested with *Nco*I and *Acc*65I and ligated into PEC610 vector also digested with the same enzymes.

PEC2122: (AD-ΔN_{28}SDF2L1ΔMIR2) Two fragments coding for the N-terminal (residues 28-94) and C-terminal (residues 150-221) were amplified by PCR using oligonucleotides #535 and #595, and #557 and #596 respectively using PEC2084 DNA as template. The oligonucleotides #595 and #596 also encoded a *Bgl*II site. The PCR products were digested with *Bgl*II and ligated. The ligated product was used as a template to amplify ΔN_{28}SDF2L1ΔMIR2 by PCR using oligonucleotides #535 and #557. The resultant PCR product was digested with *Eco*RI and *Xho*I and ligated into pGADT7-Rec plasmid also digested with the same enzymes.

PEC2123: (DBD-proHNP3Δ35-49) The proHNP3Δ35-49 DNA was amplified by PCR using oligonucleotides # 570 and #597 and plasmid PEC2119 (proHNP3Δ35-49/pCR2.1 Topo) as template. The PCR product was digested with *Eco*RI and *Bam*HI and the proHNP3Δ35-49 fragment was ligated into pGBKT7 vector also digested with the same enzymes.

PEC2124: (DBD-proRTD1aΔRLL) The proRTD1aΔRLL was amplified using oligonucleotides #512 and #547 and plasmid PEC2038 as template by PCR. The PCR product was digested using *Bam*HI and *Nco*I and ligated into pGBKT7 vector also digested with the same enzymes.

PEC2125: (DBD-RTD1a) The oligonucleotides #548 and #549 which encode the RTD1a nonamer sequence were annealed by first boiling in water bath and allowing it to cool to

room temperature. The product was then ligated into pGBKT7 vector digested with *Eco*RI and *Bam*HI.

PEC2126: (DBD-HNP3) The mature HNP3 sequence was amplified by PCR using oligonucleotides #588 and #597 and PEC2080 as template. The PCR product was digested with *Eco*RI and *Xho*I and then ligated into pGBKT7 vector also digested with *Eco*RI and *Xho*I.

PEC2127: (AD-MIR3) Oligonucleotides #557 and #600 were used to amplify the C-terminal domain of SDF2L1 (residues 151 to 221) by PCR using PEC2069 as template. The PCR product was digested with *Eco*RI and *Xho*I and ligated into PEC2084 (pGADT7-Rec) plasmid also digested with the same enzymes.

List of primers:

#506: 5'-GCGGATCCTTATAACAAACGGCAGAATCC
#512: 5'-GCCCATGGAGGCACGTCAGGCACGTGCTGATG
#513: 5'-GCGGTACCGCGGCCGCTTATAACAAACGGCAGAATCCGCGTGTGCAAATGC ACCGC
#535: 5'-CTATTCGATGATGAAGATACCCCACCAAACCC
#536: 5'-GTGAACTTGCGGGGTTTTTCAGTATCTACG
#547: 5'-GCGGATCCTTAGCAGAATCCTCGTGTGC
#548: 5'-AATTCAGGTGCATTTGCACACGAGGATTCTGCTAAG
#549: 5'-GATCCTTAGCAGAATCCTCGTGTGCAAATGCACCTG
#552: 5'-GCGAATTCGAGGCACGTCAGGCAAGAGC
#553: 5'-GCGGATCCTACAAGCCTTTCGCTGACTCTG
#556: 5'-GCTCATGAGCGAATTCGCCAAGACTGGCGCGGGGCTCG
#557: 5'- GCGTCGACTCGAGGTACCTCAGAGTTCATCGTGACCTGC
#570: 5'-GCCCATGGAATTCGAGCCACTCCAGGCAAGAGCTG
#580: 5'-GCTCATGAGCGAATTCTCGGAAGGCGGGTGCCCGCGTGGG
#581: 5'-GCGTCGACTCGAGGTACCTAAGAGCAGCGCACTGTCCATAG
#584: 5'-GCCCATGGGAGAATTCGAGTCACTCCAGGAAAGAGC
#585: 5'-GCGGATCCGGTACCTCGAGTCAGCGACAGCAGAGTCTGTAG
#586: 5'- GCCCATGGGAGAATTCGGTAACTTTCTCACAGGCC
#587: 5'-GCGGATCCGGTACCTCGAGTCACTTGCAGCACTTGGCC
#588: 5'-GCGAATTCGACTGCTATTGCAGAATACC
#597: 5'-GCGGATCCGGTACCTCGAGTCAGCAGCAGAATGCCCAGAGTC
#600: 5'-GCCCATGGAATTCGGACAGCACTGGGAGCGTGAG

Homology modeling of SDF2L1: The structure of the *C. elegans* protein, 1t9f.PDB was used for homology modeling of SDF2L1 using SWISS-MODEL. The pdb_extract program available at http://deposit.rcsb.org was then used to display the co-ordinates of the modeled structure from the PDB file created by SWISS-MODEL.

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data_UNNAMED
#
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_atom_site.label_atom_id
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_atom_site.label_seq_id
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_atom_site.Cartn_x
_atom_site.Cartn_y
_atom_site.Cartn_z
_atom_site.occupancy
_atom_site.B_iso_or_equiv
_atom_site.pdbx_PDB_model_num
ATOM   1    N   N   . LYS LYS . . 5   5   . 4.336   39.813 -22.052 1.00 50.00 .
ATOM   2    CA  CA  . LYS LYS . . 5   5   . 4.299   38.539 -21.301 1.00 50.00 .
ATOM   3    C   C   . LYS LYS . . 5   5   . 4.750   37.370 -22.197 1.00 50.00 .
ATOM   4    O   O   . LYS LYS . . 5   5   . 3.957   36.540 -22.647 1.00 50.00 .
ATOM   5    CB  CB  . LYS LYS . . 5   5   . 2.875   38.261 -20.799 1.00 50.00 .
ATOM   6    CG  CG  . LYS LYS . . 5   5   . 2.287   39.378 -19.941 1.00 50.00 .
ATOM   7    CD  CD  . LYS LYS . . 5   5   . 0.856   39.019 -19.532 1.00 50.00 .
ATOM   8    CE  CE  . LYS LYS . . 5   5   . 0.338   40.013 -18.489 1.00 50.00 .
ATOM   9    NZ  NZ  . LYS LYS . . 5   5   . -0.830  39.485 -17.770 1.00 50.00 .
ATOM   10   N   N   . THR THR . . 6   6   . 6.054   37.339 -22.472 1.00 50.00 .
ATOM   11   CA  CA  . THR THR . . 6   6   . 6.654   36.306 -23.346 1.00 50.00 .
ATOM   12   C   C   . THR THR . . 6   6   . 6.165   34.924 -22.897 1.00 50.00 .
ATOM   13   O   O   . THR THR . . 6   6   . 6.216   34.614 -21.707 1.00 50.00 .
ATOM   14   CB  CB  . THR THR . . 6   6   . 8.188   36.339 -23.278 1.00 50.00 .
ATOM   15   OG1 OG1 . THR THR . . 6   6   . 8.616   36.094 -21.935 1.00 50.00 .
ATOM   16   CG2 CG2 . THR THR . . 6   6   . 8.749   37.667 -23.797 1.00 50.00 .
ATOM   17   N   N   . GLY GLY . . 7   7   . 5.707   34.130 -23.868 1.00 50.00 .
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ATOM   20   O   O   . GLY GLY . . 7   7   . 6.004   30.584 -23.034 1.00 50.00 .
ATOM   21   N   N   . ALA ALA . . 8   8   . 7.447   32.298 -22.863 1.00 50.00 .
ATOM   22   CA  CA  . ALA ALA . . 8   8   . 8.611   31.518 -22.416 1.00 50.00 .
ATOM   23   C   C   . ALA ALA . . 8   8   . 9.242   32.098 -21.137 1.00 50.00 .
ATOM   24   O   O   . ALA ALA . . 8   8   . 10.401  31.844 -20.819 1.00 50.00 .
ATOM   25   CB  CB  . ALA ALA . . 8   8   . 9.617   31.470 -23.572 1.00 50.00 .
ATOM   26   N   N   . GLU GLU . . 9   9   . 8.419   32.768 -20.333 1.00 50.00 .
ATOM   27   CA  CA  . GLU GLU . . 9   9   . 8.857   33.302 -19.024 1.00 50.00 .
ATOM   28   C   C   . GLU GLU . . 9   9   . 8.452   32.363 -17.874 1.00 50.00 .
ATOM   29   O   O   . GLU GLU . . 9   9   . 7.999   32.765 -16.802 1.00 50.00 .
ATOM   30   CB  CB  . GLU GLU . . 9   9   . 8.314   34.718 -18.822 1.00 50.00 .
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ATOM   31   CG  CG  . GLU GLU . . 9   9   . 6.786   34.757 -18.832 1.00 50.00 .
ATOM   32   CD  CD  . GLU GLU . . 9   9   . 6.319   36.182 -18.611 1.00 50.00 .
ATOM   33   OE1 OE1 . GLU GLU . . 9   9   . 6.521   37.014 -19.526 1.00 50.00 .
ATOM   34   OE2 OE2 . GLU GLU . . 9   9   . 5.851   36.409 -17.488 1.00 50.00 .
ATOM   35   N   N   . LEU LEU . . 10  10  . 8.484   31.083 -18.211 1.00 50.00 .
ATOM   36   CA  CA  . LEU LEU . . 10  10  . 8.123   29.997 -17.299 1.00 50.00 .
ATOM   37   C   C   . LEU LEU . . 10  10  . 8.991   29.932 -16.059 1.00 50.00 .
ATOM   38   O   O   . LEU LEU . . 10  10  . 10.160  30.303 -16.069 1.00 50.00 .
ATOM   39   CB  CB  . LEU LEU . . 10  10  . 8.241   28.658 -18.024 1.00 50.00 .
ATOM   40   CG  CG  . LEU LEU . . 10  10  . 7.269   28.595 -19.188 1.00 50.00 .
ATOM   41   CD1 CD1 . LEU LEU . . 10  10  . 7.436   27.269 -19.920 1.00 50.00 .
ATOM   42   CD2 CD2 . LEU LEU . . 10  10  . 5.874   28.797 -18.631 1.00 50.00 .
ATOM   43   N   N   . VAL VAL . . 11  11  . 8.355   29.435 -15.018 1.00 50.00 .
ATOM   44   CA  CA  . VAL VAL . . 11  11  . 9.020   29.119 -13.758 1.00 50.00 .
ATOM   45   C   C   . VAL VAL . . 11  11  . 9.818   27.822 -13.987 1.00 50.00 .
ATOM   46   O   O   . VAL VAL . . 11  11  . 9.336   26.867 -14.597 1.00 50.00 .
ATOM   47   CB  CB  . VAL VAL . . 11  11  . 7.947   28.960 -12.678 1.00 50.00 .
ATOM   48   CG1 CG1 . VAL VAL . . 11  11  . 8.652   28.793 -11.353 1.00 50.00 .
ATOM   49   CG2 CG2 . VAL VAL . . 11  11  . 7.039   30.186 -12.533 1.00 50.00 .
ATOM   50   N   N   . THR THR . . 12  12  . 11.032  27.808 -13.457 1.00 50.00 .
ATOM   51   CA  CA  . THR THR . . 12  12  . 11.964  26.691 -13.687 1.00 50.00 .
ATOM   52   C   C   . THR THR . . 12  12  . 12.297  25.966 -12.379 1.00 50.00 .
ATOM   53   O   O   . THR THR . . 12  12  . 12.201  26.519 -11.289 1.00 50.00 .
ATOM   54   CB  CB  . THR THR . . 12  12  . 13.249  27.176 -14.371 1.00 50.00 .
ATOM   55   OG1 OG1 . THR THR . . 12  12  . 13.946  28.081 -13.510 1.00 50.00 .
ATOM   56   CG2 CG2 . THR THR . . 12  12  . 12.957  27.827 -15.730 1.00 50.00 .
ATOM   57   N   N   . CYS CYS . . 13  13  . 12.631  24.692 -12.505 1.00 50.00 .
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ATOM   61   CB  CB  . CYS CYS . . 13  13  . 13.201  22.412 -11.828 1.00 50.00 .
ATOM   62   SG  SG  . CYS CYS . . 13  13  . 12.736  21.208 -10.535 1.00 50.00 .
ATOM   63   N   N   . GLY GLY . . 14  14  . 14.264  24.340 -9.377  1.00 50.00 .
ATOM   64   CA  CA  . GLY GLY . . 14  14  . 15.394  24.823 -8.562  1.00 50.00 .
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ATOM   75   C   C   . VAL VAL . . 16  16  . 13.486  31.994 -6.291  1.00 50.00 .
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ATOM   77   CB  CB  . VAL VAL . . 16  16  . 15.584  31.129 -5.235  1.00 50.00 .
ATOM   78   CG1 CG1 . VAL VAL . . 16  16  . 15.266  31.698 -3.848  1.00 50.00 .
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ATOM   82   C   C   . LEU LEU . . 17  17  . 10.507  33.348 -4.492  1.00 50.00 .
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ATOM   85   CG  CG  . LEU LEU . . 17  17  . 10.214  31.530 -7.260  1.00 50.00 .
ATOM   86   CD1 CD1 . LEU LEU . . 17  17  . 9.112   31.092 -6.302  1.00 50.00 .
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ATOM   91   O   O   . LYS LYS . . 18  18  . 7.048   34.562 -5.305  1.00 50.00 .
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ATOM   104  CD2 CD2 . LEU LEU . . 19  19  . 4.508   29.850 -3.444  1.00 50.00 .
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ATOM   106  CA  CA  . LEU LEU . . 20  20  . 2.723   36.720 -3.292  1.00 50.00 .
ATOM   107  C   C   . LEU LEU . . 20  20  . 1.264   36.327 -3.411  1.00 50.00 .
ATOM   108  O   O   . LEU LEU . . 20  20  . 0.825   35.890 -4.471  1.00 50.00 .
ATOM   109  CB  CB  . LEU LEU . . 20  20  . 2.932   37.905 -4.238  1.00 50.00 .
ATOM   110  CG  CG  . LEU LEU . . 20  20  . 4.319   38.510 -4.129  1.00 50.00 .
ATOM   111  CD1 CD1 . LEU LEU . . 20  20  . 4.443   39.755 -4.998  1.00 50.00 .
ATOM   112  CD2 CD2 . LEU LEU . . 20  20  . 4.396   39.015 -2.721  1.00 50.00 .
ATOM   113  N   N   . ASN ASN . . 21  21  . 0.530   36.499 -2.322  1.00 50.00 .
ATOM   114  CA  CA  . ASN ASN . . 21  21  . -0.925  36.292 -2.350  1.00 50.00 .
ATOM   115  C   C   . ASN ASN . . 21  21  . -1.549  37.414 -3.184  1.00 50.00 .
ATOM   116  O   O   . ASN ASN . . 21  21  . -1.486  38.582 -2.818  1.00 50.00 .
ATOM   117  CB  CB  . ASN ASN . . 21  21  . -1.486  36.292 -0.931  1.00 50.00 .
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ATOM   119  OD1 OD1 . ASN ASN . . 21  21  . -3.864  36.606 -1.154  1.00 50.00 .
ATOM   120  ND2 ND2 . ASN ASN . . 21  21  . -3.110  34.532 -0.815  1.00 50.00 .
ATOM   121  N   N   . THR THR . . 22  22  . -2.249  37.020 -4.234  1.00 50.00 .
ATOM   122  CA  CA  . THR THR . . 22  22  . -2.805  37.970 -5.225  1.00 50.00 .
ATOM   123  C   C   . THR THR . . 22  22  . -3.732  39.056 -4.661  1.00 50.00 .
ATOM   124  O   O   . THR THR . . 22  22  . -3.582  40.227 -4.994  1.00 50.00 .
ATOM   125  CB  CB  . THR THR . . 22  22  . -3.566  37.217 -6.310  1.00 50.00 .
ATOM   126  OG1 OG1 . THR THR . . 22  22  . -4.619  36.464 -5.693  1.00 50.00 .
ATOM   127  CG2 CG2 . THR THR . . 22  22  . -2.589  36.376 -7.135  1.00 50.00 .
ATOM   128  N   N   . HIS HIS . . 23  23  . -4.609  38.659 -3.741  1.00 50.00 .
ATOM   129  CA  CA  . HIS HIS . . 23  23  . -5.629  39.565 -3.182  1.00 50.00 .
ATOM   130  C   C   . HIS HIS . . 23  23  . -5.125  40.378 -1.984  1.00 50.00 .
ATOM   131  O   O   . HIS HIS . . 23  23  . -5.618  41.471 -1.721  1.00 50.00 .
ATOM   132  CB  CB  . HIS HIS . . 23  23  . -6.886  38.782 -2.792  1.00 50.00 .
ATOM   133  CG  CG  . HIS HIS . . 23  23  . -7.586  38.191 -4.019  1.00 50.00 .
ATOM   134  ND1 ND1 . HIS HIS . . 23  23  . -7.338  37.002 -4.557  1.00 50.00 .
ATOM   135  CD2 CD2 . HIS HIS . . 23  23  . -8.584  38.746 -4.701  1.00 50.00 .
ATOM   136  CE1 CE1 . HIS HIS . . 23  23  . -8.181  36.817 -5.566  1.00 50.00 .
ATOM   137  NE2 NE2 . HIS HIS . . 23  23  . -8.955  37.895 -5.655  1.00 50.00 .
ATOM   138  N   N   . HIS HIS . . 24  24  . -4.146  39.824 -1.273  1.00 50.00 .
ATOM   139  CA  CA  . HIS HIS . . 24  24  . -3.619  40.470 -0.062  1.00 50.00 .
ATOM   140  C   C   . HIS HIS . . 24  24  . -2.282  41.179 -0.241  1.00 50.00 .
ATOM   141  O   O   . HIS HIS . . 24  24  . -2.010  42.086 0.539   1.00 50.00 .
ATOM   142  CB  CB  . HIS HIS . . 24  24  . -3.518  39.494 1.105   1.00 50.00 .
ATOM   143  CG  CG  . HIS HIS . . 24  24  . -4.922  39.095 1.529   1.00 50.00 .
ATOM   144  ND1 ND1 . HIS HIS . . 24  24  . -5.790  39.848 2.197   1.00 50.00 .
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ATOM   145  CD2 CD2 . HIS HIS . . 24  24  . -5.426  37.883 1.398   1.00 50.00 .
ATOM   146  CE1 CE1 . HIS HIS . . 24  24  . -6.830  39.091 2.514   1.00 50.00 .
ATOM   147  NE2 NE2 . HIS HIS . . 24  24  . -6.588  37.881 2.030   1.00 50.00 .
ATOM   148  N   N   . ARG ARG . . 25  25  . -1.457  40.683 -1.173  1.00 50.00 .
ATOM   149  CA  CA  . ARG ARG . . 25  25  . -0.115  41.209 -1.531  1.00 50.00 .
ATOM   150  C   C   . ARG ARG . . 25  25  . 1.005   40.687 -0.610  1.00 50.00 .
ATOM   151  O   O   . ARG ARG . . 25  25  . 2.200   40.870 -0.838  1.00 50.00 .
ATOM   152  CB  CB  . ARG ARG . . 25  25  . -0.158  42.742 -1.571  1.00 50.00 .
ATOM   153  CG  CG  . ARG ARG . . 25  25  . 1.169   43.466 -1.775  1.00 50.00 .
ATOM   154  CD  CD  . ARG ARG . . 25  25  . 0.962   44.961 -1.542  1.00 50.00 .
ATOM   155  NE  NE  . ARG ARG . . 25  25  . 0.057   45.496 -2.575  1.00 50.00 .
ATOM   156  CZ  CZ  . ARG ARG . . 25  25  . 0.387   45.766 -3.841  1.00 50.00 .
ATOM   157  NH1 NH1 . ARG ARG . . 25  25  . 1.622   45.573 -4.287  1.00 50.00 .
ATOM   158  NH2 NH2 . ARG ARG . . 25  25  . -0.529  46.230 -4.676  1.00 50.00 .
ATOM   159  N   N   . VAL VAL . . 26  26  . 0.580   39.959 0.407   1.00 50.00 .
ATOM   160  CA  CA  . VAL VAL . . 26  26  . 1.457   39.338 1.395   1.00 50.00 .
ATOM   161  C   C   . VAL VAL . . 26  26  . 2.332   38.251 0.738   1.00 50.00 .
ATOM   162  O   O   . VAL VAL . . 26  26  . 1.868   37.544 -0.156  1.00 50.00 .
ATOM   163  CB  CB  . VAL VAL . . 26  26  . 0.467   38.841 2.451   1.00 50.00 .
ATOM   164  CG1 CG1 . VAL VAL . . 26  26  . -0.273  37.534 2.162   1.00 50.00 .
ATOM   165  CG2 CG2 . VAL VAL . . 26  26  . 1.170   38.740 3.768   1.00 50.00 .
ATOM   166  N   N   . ARG ARG . . 27  27  . 3.528   38.052 1.275   1.00 50.00 .
ATOM   167  CA  CA  . ARG ARG . . 27  27  . 4.465   37.037 0.754   1.00 50.00 .
ATOM   168  C   C   . ARG ARG . . 27  27  . 4.436   35.818 1.674   1.00 50.00 .
ATOM   169  O   O   . ARG ARG . . 27  27  . 4.551   35.974 2.889   1.00 50.00 .
ATOM   170  CB  CB  . ARG ARG . . 27  27  . 5.919   37.528 0.779   1.00 50.00 .
ATOM   171  CG  CG  . ARG ARG . . 27  27  . 6.279   38.707 -0.124  1.00 50.00 .
ATOM   172  CD  CD  . ARG ARG . . 27  27  . 5.676   40.091 0.215   1.00 50.00 .
ATOM   173  NE  NE  . ARG ARG . . 27  27  . 5.645   41.023 -0.948  1.00 50.00 .
ATOM   174  CZ  CZ  . ARG ARG . . 27  27  . 6.652   41.340 -1.778  1.00 50.00 .
ATOM   175  NH1 NH1 . ARG ARG . . 27  27  . 7.846   40.782 -1.643  1.00 50.00 .
ATOM   176  NH2 NH2 . ARG ARG . . 27  27  . 6.458   42.147 -2.814  1.00 50.00 .
ATOM   177  N   N   . LEU LEU . . 28  28  . 4.441   34.640 1.062   1.00 50.00 .
ATOM   178  CA  CA  . LEU LEU . . 28  28  . 4.605   33.369 1.790   1.00 50.00 .
ATOM   179  C   C   . LEU LEU . . 28  28  . 5.946   33.418 2.536   1.00 50.00 .
ATOM   180  O   O   . LEU LEU . . 28  28  . 6.969   33.803 1.971   1.00 50.00 .
ATOM   181  CB  CB  . LEU LEU . . 28  28  . 4.584   32.194 0.806   1.00 50.00 .
ATOM   182  CG  CG  . LEU LEU . . 28  28  . 4.492   30.839 1.518   1.00 50.00 .
ATOM   183  CD1 CD1 . LEU LEU . . 28  28  . 3.175   30.710 2.289   1.00 50.00 .
ATOM   184  CD2 CD2 . LEU LEU . . 28  28  . 4.621   29.702 0.505   1.00 50.00 .
ATOM   185  N   N   . HIS HIS . . 29  29  . 5.887   33.068 3.811   1.00 50.00 .
ATOM   186  CA  CA  . HIS HIS . . 29  29  . 6.997   33.303 4.744   1.00 50.00 .
ATOM   187  C   C   . HIS HIS . . 29  29  . 7.055   32.216 5.818   1.00 50.00 .
ATOM   188  O   O   . HIS HIS . . 29  29  . 6.068   31.536 6.105   1.00 50.00 .
ATOM   189  CB  CB  . HIS HIS . . 29  29  . 6.766   34.673 5.387   1.00 50.00 .
ATOM   190  CG  CG  . HIS HIS . . 29  29  . 7.894   35.098 6.327   1.00 50.00 .
ATOM   191  ND1 ND1 . HIS HIS . . 29  29  . 7.951   34.840 7.629   1.00 50.00 .
ATOM   192  CD2 CD2 . HIS HIS . . 29  29  . 8.901   35.907 6.026   1.00 50.00 .
ATOM   193  CE1 CE1 . HIS HIS . . 29  29  . 8.993   35.493 8.134   1.00 50.00 .
ATOM   194  NE2 NE2 . HIS HIS . . 29  29  . 9.571   36.165 7.142   1.00 50.00 .
ATOM   195  N   N   . SER SER . . 30  30  . 8.275   31.997 6.280   1.00 50.00 .
ATOM   196  CA  CA  . SER SER . . 30  30  . 8.550   31.090 7.401   1.00 50.00 .
ATOM   197  C   C   . SER SER . . 30  30  . 9.743   31.636 8.197   1.00 50.00 .
ATOM   198  O   O   . SER SER . . 30  30  . 10.395  32.608 7.817   1.00 50.00 .
ATOM   199  CB  CB  . SER SER . . 30  30  . 8.802   29.676 6.872   1.00 50.00 .
ATOM   200  OG  OG  . SER SER . . 30  30  . 8.940   28.766 7.962   1.00 50.00 .
ATOM   201  N   N   . HIS HIS . . 31  31  . 9.992   31.004 9.333   1.00 50.00 .
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ATOM   202  CA  CA  . HIS HIS . . 31  31  . 10.960  31.471 10.346  1.00 50.00 .
ATOM   203  C   C   . HIS HIS . . 31  31  . 11.157  30.403 11.427  1.00 50.00 .
ATOM   204  O   O   . HIS HIS . . 31  31  . 10.459  29.396 11.442  1.00 50.00 .
ATOM   205  CB  CB  . HIS HIS . . 31  31  . 10.463  32.796 10.967  1.00 50.00 .
ATOM   206  CG  CG  . HIS HIS . . 31  31  . 8.957   32.851 11.288  1.00 50.00 .
ATOM   207  ND1 ND1 . HIS HIS . . 31  31  . 8.209   33.948 11.239  1.00 50.00 .
ATOM   208  CD2 CD2 . HIS HIS . . 31  31  . 8.133   31.816 11.454  1.00 50.00 .
ATOM   209  CE1 CE1 . HIS HIS . . 31  31  . 6.931   33.575 11.307  1.00 50.00 .
ATOM   210  NE2 NE2 . HIS HIS . . 31  31  . 6.886   32.253 11.446  1.00 50.00 .
ATOM   211  N   N   . ASP ASP . . 32  32  . 11.970  30.717 12.428  1.00 50.00 .
ATOM   212  CA  CA  . ASP ASP . . 32  32  . 12.251  29.775 13.528  1.00 50.00 .
ATOM   213  C   C   . ASP ASP . . 32  32  . 11.218  29.890 14.670  1.00 50.00 .
ATOM   214  O   O   . ASP ASP . . 32  32  . 11.548  29.975 15.851  1.00 50.00 .
ATOM   215  CB  CB  . ASP ASP . . 32  32  . 13.688  29.986 14.019  1.00 50.00 .
ATOM   216  CG  CG  . ASP ASP . . 32  32  . 14.204  28.804 14.853  1.00 50.00 .
ATOM   217  OD1 OD1 . ASP ASP . . 32  32  . 13.388  28.104 15.493  1.00 50.00 .
ATOM   218  OD2 OD2 . ASP ASP . . 32  32  . 15.425  28.567 14.760  1.00 50.00 .
ATOM   219  N   N   . ILE ILE . . 33  33  . 9.959   30.006 14.272  1.00 50.00 .
ATOM   220  CA  CA  . ILE ILE . . 33  33  . 8.827   30.078 15.212  1.00 50.00 .
ATOM   221  C   C   . ILE ILE . . 33  33  . 7.935   28.869 14.937  1.00 50.00 .
ATOM   222  O   O   . ILE ILE . . 33  33  . 7.428   28.696 13.834  1.00 50.00 .
ATOM   223  CB  CB  . ILE ILE . . 33  33  . 8.078   31.411 15.053  1.00 50.00 .
ATOM   224  CG1 CG1 . ILE ILE . . 33  33  . 9.080   32.535 15.353  1.00 50.00 .
ATOM   225  CG2 CG2 . ILE ILE . . 33  33  . 6.841   31.480 15.969  1.00 50.00 .
ATOM   226  CD1 CD1 . ILE ILE . . 33  33  . 8.609   33.954 15.068  1.00 50.00 .
ATOM   227  N   N   . LYS LYS . . 34  34  . 7.775   28.052 15.965  1.00 50.00 .
ATOM   228  CA  CA  . LYS LYS . . 34  34  . 6.979   26.820 15.846  1.00 50.00 .
ATOM   229  C   C   . LYS LYS . . 34  34  . 5.555   27.022 16.371  1.00 50.00 .
ATOM   230  O   O   . LYS LYS . . 34  34  . 5.292   27.903 17.190  1.00 50.00 .
ATOM   231  CB  CB  . LYS LYS . . 34  34  . 7.654   25.665 16.588  1.00 50.00 .
ATOM   232  CG  CG  . LYS LYS . . 34  34  . 9.017   25.333 15.976  1.00 50.00 .
ATOM   233  CD  CD  . LYS LYS . . 34  34  . 9.539   24.001 16.508  1.00 50.00 .
ATOM   234  CE  CE  . LYS LYS . . 34  34  . 10.871  23.653 15.844  1.00 50.00 .
ATOM   235  NZ  NZ  . LYS LYS . . 34  34  . 11.269  22.276 16.167  1.00 50.00 .
ATOM   236  N   N   . TYR TYR . . 35  35  . 4.642   26.249 15.794  1.00 50.00 .
ATOM   237  CA  CA  . TYR TYR . . 35  35  . 3.239   26.212 16.244  1.00 50.00 .
ATOM   238  C   C   . TYR TYR . . 35  35  . 3.149   25.698 17.687  1.00 50.00 .
ATOM   239  O   O   . TYR TYR . . 35  35  . 3.965   24.892 18.128  1.00 50.00 .
ATOM   240  CB  CB  . TYR TYR . . 35  35  . 2.391   25.310 15.340  1.00 50.00 .
ATOM   241  CG  CG  . TYR TYR . . 35  35  . 2.171   25.894 13.944  1.00 50.00 .
ATOM   242  CD1 CD1 . TYR TYR . . 35  35  . 1.526   27.117 13.788  1.00 50.00 .
ATOM   243  CD2 CD2 . TYR TYR . . 35  35  . 2.494   25.132 12.828  1.00 50.00 .
ATOM   244  CE1 CE1 . TYR TYR . . 35  35  . 1.199   27.572 12.520  1.00 50.00 .
ATOM   245  CE2 CE2 . TYR TYR . . 35  35  . 2.164   25.585 11.558  1.00 50.00 .
ATOM   246  CZ  CZ  . TYR TYR . . 35  35  . 1.519   26.807 11.410  1.00 50.00 .
ATOM   247  OH  OH  . TYR TYR . . 35  35  . 1.271   27.326 10.186  1.00 50.00 .
ATOM   248  N   N   . GLY GLY . . 36  36  . 2.123   26.194 18.380  1.00 50.00 .
ATOM   249  CA  CA  . GLY GLY . . 36  36  . 1.807   25.751 19.751  1.00 50.00 .
ATOM   250  C   C   . GLY GLY . . 36  36  . 0.377   25.204 19.815  1.00 50.00 .
ATOM   251  O   O   . GLY GLY . . 36  36  . -0.435  25.697 20.593  1.00 50.00 .
ATOM   252  N   N   . SER SER . . 37  37  . 0.072   24.283 18.901  1.00 50.00 .
ATOM   253  CA  CA  . SER SER . . 37  37  . -1.300  23.763 18.719  1.00 50.00 .
ATOM   254  C   C   . SER SER . . 37  37  . -1.365  22.415 17.985  1.00 50.00 .
ATOM   255  O   O   . SER SER . . 37  37  . -1.501  21.381 18.631  1.00 50.00 .
ATOM   256  CB  CB  . SER SER . . 37  37  . -2.192  24.810 18.028  1.00 50.00 .
ATOM   257  OG  OG  . SER SER . . 37  37  . -1.620  25.217 16.775  1.00 50.00 .
ATOM   258  N   N   . GLY GLY . . 38  38  . -1.220  22.462 16.653  1.00 50.00 .
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ATOM   259  CA  CA  . GLY GLY . . 38  38  . -1.400  21.294 15.769  1.00 50.00 .
ATOM   260  C   C   . GLY GLY . . 38  38  . -0.297  20.249 15.954  1.00 50.00 .
ATOM   261  O   O   . GLY GLY . . 38  38  . -0.228  19.540 16.953  1.00 50.00 .
ATOM   262  N   N   . SER SER . . 39  39  . 0.612   20.246 14.990  1.00 50.00 .
ATOM   263  CA  CA  . SER SER . . 39  39  . 1.776   19.337 15.008  1.00 50.00 .
ATOM   264  C   C   . SER SER . . 39  39  . 2.931   19.871 15.872  1.00 50.00 .
ATOM   265  O   O   . SER SER . . 39  39  . 3.901   19.164 16.133  1.00 50.00 .
ATOM   266  CB  CB  . SER SER . . 39  39  . 2.264   19.098 13.579  1.00 50.00 .
ATOM   267  OG  OG  . SER SER . . 39  39  . 2.718   20.334 13.015  1.00 50.00 .
ATOM   268  N   N   . GLY GLY . . 40  40  . 2.853   21.179 16.165  1.00 50.00 .
ATOM   269  CA  CA  . GLY GLY . . 40  40  . 3.920   21.931 16.854  1.00 50.00 .
ATOM   270  C   C   . GLY GLY . . 40  40  . 5.194   22.039 16.001  1.00 50.00 .
ATOM   271  O   O   . GLY GLY . . 40  40  . 6.291   22.241 16.514  1.00 50.00 .
ATOM   272  N   N   . GLN GLN . . 41  41  . 5.019   21.887 14.690  1.00 50.00 .
ATOM   273  CA  CA  . GLN GLN . . 41  41  . 6.103   22.081 13.712  1.00 50.00 .
ATOM   274  C   C   . GLN GLN . . 41  41  . 6.284   23.577 13.435  1.00 50.00 .
ATOM   275  O   O   . GLN GLN . . 41  41  . 5.470   24.404 13.851  1.00 50.00 .
ATOM   276  CB  CB  . GLN GLN . . 41  41  . 5.789   21.339 12.414  1.00 50.00 .
ATOM   277  CG  CG  . GLN GLN . . 41  41  . 5.767   19.823 12.622  1.00 50.00 .
ATOM   278  CD  CD  . GLN GLN . . 41  41  . 5.399   19.092 11.333  1.00 50.00 .
ATOM   279  OE1 OE1 . GLN GLN . . 41  41  . 4.711   19.601 10.456  1.00 50.00 .
ATOM   280  NE2 NE2 . GLN GLN . . 41  41  . 5.854   17.865 11.246  1.00 50.00 .
ATOM   281  N   N   . GLN GLN . . 42  42  . 7.344   23.892 12.700  1.00 50.00 .
ATOM   282  CA  CA  . GLN GLN . . 42  42  . 7.651   25.277 12.304  1.00 50.00 .
ATOM   283  C   C   . GLN GLN . . 42  42  . 6.478   25.875 11.509  1.00 50.00 .
ATOM   284  O   O   . GLN GLN . . 42  42  . 5.955   25.248 10.585  1.00 50.00 .
ATOM   285  CB  CB  . GLN GLN . . 42  42  . 8.920   25.242 11.456  1.00 50.00 .
ATOM   286  CG  CG  . GLN GLN . . 42  42  . 9.473   26.646 11.221  1.00 50.00 .
ATOM   287  CD  CD  . GLN GLN . . 42  42  . 10.772  26.607 10.412  1.00 50.00 .
ATOM   288  OE1 OE1 . GLN GLN . . 42  42  . 10.855  26.933 9.232   1.00 50.00 .
ATOM   289  NE2 NE2 . GLN GLN . . 42  42  . 11.837  26.191 11.063  1.00 50.00 .
ATOM   290  N   N   . SER SER . . 43  43  . 6.022   27.030 11.983  1.00 50.00 .
ATOM   291  CA  CA  . SER SER . . 43  43  . 4.886   27.753 11.382  1.00 50.00 .
ATOM   292  C   C   . SER SER . . 43  43  . 5.217   28.355 10.009  1.00 50.00 .
ATOM   293  O   O   . SER SER . . 43  43  . 6.379   28.551 9.655   1.00 50.00 .
ATOM   294  CB  CB  . SER SER . . 43  43  . 4.394   28.864 12.314  1.00 50.00 .
ATOM   295  OG  OG  . SER SER . . 43  43  . 5.343   29.929 12.408  1.00 50.00 .
ATOM   296  N   N   . VAL VAL . . 44  44  . 4.153   28.556 9.240   1.00 50.00 .
ATOM   297  CA  CA  . VAL VAL . . 44  44  . 4.206   29.235 7.930   1.00 50.00 .
ATOM   298  C   C   . VAL VAL . . 44  44  . 3.184   30.377 7.966   1.00 50.00 .
ATOM   299  O   O   . VAL VAL . . 44  44  . 2.032   30.189 8.363   1.00 50.00 .
ATOM   300  CB  CB  . VAL VAL . . 44  44  . 3.871   28.252 6.792   1.00 50.00 .
ATOM   301  CG1 CG1 . VAL VAL . . 44  44  . 3.835   28.932 5.420   1.00 50.00 .
ATOM   302  CG2 CG2 . VAL VAL . . 44  44  . 4.917   27.143 6.709   1.00 50.00 .
ATOM   303  N   N   . THR THR . . 45  45  . 3.637   31.546 7.542   1.00 50.00 .
ATOM   304  CA  CA  . THR THR . . 45  45  . 2.775   32.738 7.498   1.00 50.00 .
ATOM   305  C   C   . THR THR . . 45  45  . 2.911   33.439 6.139   1.00 50.00 .
ATOM   306  O   O   . THR THR . . 45  45  . 3.518   32.930 5.196   1.00 50.00 .
ATOM   307  CB  CB  . THR THR . . 45  45  . 3.130   33.721 8.630   1.00 50.00 .
ATOM   308  OG1 OG1 . THR THR . . 45  45  . 4.403   34.303 8.346   1.00 50.00 .
ATOM   309  CG2 CG2 . THR THR . . 45  45  . 3.110   33.073 10.020  1.00 50.00 .
ATOM   310  N   N   . GLY GLY . . 46  46  . 2.187   34.551 6.048   1.00 50.00 .
ATOM   311  CA  CA  . GLY GLY . . 46  46  . 2.352   35.515 4.960   1.00 50.00 .
ATOM   312  C   C   . GLY GLY . . 46  46  . 2.692   36.841 5.639   1.00 50.00 .
ATOM   313  O   O   . GLY GLY . . 46  46  . 1.915   37.306 6.476   1.00 50.00 .
ATOM   314  N   N   . VAL VAL . . 47  47  . 3.677   37.531 5.077   1.00 50.00 .
ATOM   315  CA  CA  . VAL VAL . . 47  47  . 4.118   38.851 5.589   1.00 50.00 .
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ATOM   316  C   C   . VAL VAL . . 47  47  . 3.891   40.014 4.607   1.00 50.00 .
ATOM   317  O   O   . VAL VAL . . 47  47  . 4.162   39.903 3.409   1.00 50.00 .
ATOM   318  CB  CB  . VAL VAL . . 47  47  . 5.591   38.804 6.019   1.00 50.00 .
ATOM   319  CG1 CG1 . VAL VAL . . 47  47  . 5.767   37.833 7.189   1.00 50.00 .
ATOM   320  CG2 CG2 . VAL VAL . . 47  47  . 6.537   38.449 4.862   1.00 50.00 .
ATOM   321  N   N   . GLU GLU . . 48  48  . 3.391   41.119 5.154   1.00 50.00 .
ATOM   322  CA  CA  . GLU GLU . . 48  48  . 3.235   42.389 4.419   1.00 50.00 .
ATOM   323  C   C   . GLU GLU . . 48  48  . 4.522   43.217 4.592   1.00 50.00 .
ATOM   324  O   O   . GLU GLU . . 48  48  . 4.538   44.320 5.136   1.00 50.00 .
ATOM   325  CB  CB  . GLU GLU . . 48  48  . 1.988   43.116 4.947   1.00 50.00 .
ATOM   326  CG  CG  . GLU GLU . . 48  48  . 1.610   44.362 4.127   1.00 50.00 .
ATOM   327  CD  CD  . GLU GLU . . 48  48  . 1.307   44.090 2.646   1.00 50.00 .
ATOM   328  OE1 OE1 . GLU GLU . . 48  48  . 1.077   42.911 2.288   1.00 50.00 .
ATOM   329  OE2 OE2 . GLU GLU . . 48  48  . 1.361   45.077 1.882   1.00 50.00 .
ATOM   330  N   N   . ALA ALA . . 49  49  . 5.626   42.550 4.275   1.00 50.00 .
ATOM   331  CA  CA  . ALA ALA . . 49  49  . 6.981   43.098 4.444   1.00 50.00 .
ATOM   332  C   C   . ALA ALA . . 49  49  . 7.855   42.639 3.275   1.00 50.00 .
ATOM   333  O   O   . ALA ALA . . 49  49  . 8.368   41.518 3.257   1.00 50.00 .
ATOM   334  CB  CB  . ALA ALA . . 49  49  . 7.553   42.645 5.794   1.00 50.00 .
ATOM   335  N   N   . SER SER . . 50  50  . 7.866   43.485 2.245   1.00 50.00 .
ATOM   336  CA  CA  . SER SER . . 50  50  . 8.617   43.241 0.994   1.00 50.00 .
ATOM   337  C   C   . SER SER . . 50  50  . 10.072  42.841 1.273   1.00 50.00 .
ATOM   338  O   O   . SER SER . . 50  50  . 10.585  43.052 2.370   1.00 50.00 .
ATOM   339  CB  CB  . SER SER . . 50  50  . 8.618   44.498 0.115   1.00 50.00 .
ATOM   340  OG  OG  . SER SER . . 50  50  . 9.337   44.266 -1.104  1.00 50.00 .
ATOM   341  N   N   . ASP ASP . . 51  51  . 10.660  42.186 0.275   1.00 50.00 .
ATOM   342  CA  CA  . ASP ASP . . 51  51  . 12.078  41.784 0.238   1.00 50.00 .
ATOM   343  C   C   . ASP ASP . . 51  51  . 12.651  41.376 1.606   1.00 50.00 .
ATOM   344  O   O   . ASP ASP . . 51  51  . 13.165  42.186 2.375   1.00 50.00 .
ATOM   345  CB  CB  . ASP ASP . . 51  51  . 12.910  42.914 -0.390  1.00 50.00 .
ATOM   346  CG  CG  . ASP ASP . . 51  51  . 13.852  42.410 -1.488  1.00 50.00 .
ATOM   347  OD1 OD1 . ASP ASP . . 51  51  . 13.880  41.179 -1.687  1.00 50.00 .
ATOM   348  OD2 OD2 . ASP ASP . . 51  51  . 14.528  43.254 -2.107  1.00 50.00 .
ATOM   349  N   N   . ASP ASP . . 52  52  . 12.462  40.102 1.911   1.00 50.00 .
ATOM   350  CA  CA  . ASP ASP . . 52  52  . 12.995  39.503 3.143   1.00 50.00 .
ATOM   351  C   C   . ASP ASP . . 52  52  . 13.523  38.114 2.775   1.00 50.00 .
ATOM   352  O   O   . ASP ASP . . 52  52  . 12.781  37.322 2.203   1.00 50.00 .
ATOM   353  CB  CB  . ASP ASP . . 52  52  . 11.877  39.410 4.197   1.00 50.00 .
ATOM   354  CG  CG  . ASP ASP . . 52  52  . 12.363  38.970 5.587   1.00 50.00 .
ATOM   355  OD1 OD1 . ASP ASP . . 52  52  . 13.506  38.464 5.690   1.00 50.00 .
ATOM   356  OD2 OD2 . ASP ASP . . 52  52  . 11.534  38.980 6.520   1.00 50.00 .
ATOM   357  N   N   . ALA ALA . . 53  53  . 14.722  37.796 3.251   1.00 50.00 .
ATOM   358  CA  CA  . ALA ALA . . 53  53  . 15.370  36.487 3.013   1.00 50.00 .
ATOM   359  C   C   . ALA ALA . . 53  53  . 14.522  35.286 3.469   1.00 50.00 .
ATOM   360  O   O   . ALA ALA . . 53  53  . 14.436  34.284 2.762   1.00 50.00 .
ATOM   361  CB  CB  . ALA ALA . . 53  53  . 16.723  36.449 3.725   1.00 50.00 .
ATOM   362  N   N   . ASN ASN . . 54  54  . 13.787  35.470 4.567   1.00 50.00 .
ATOM   363  CA  CA  . ASN ASN . . 54  54  . 12.849  34.454 5.093   1.00 50.00 .
ATOM   364  C   C   . ASN ASN . . 54  54  . 11.577  34.253 4.245   1.00 50.00 .
ATOM   365  O   O   . ASN ASN . . 54  54  . 10.696  33.468 4.580   1.00 50.00 .
ATOM   366  CB  CB  . ASN ASN . . 54  54  . 12.460  34.810 6.527   1.00 50.00 .
ATOM   367  CG  CG  . ASN ASN . . 54  54  . 13.634  34.683 7.490   1.00 50.00 .
ATOM   368  OD1 OD1 . ASN ASN . . 54  54  . 14.278  33.654 7.604   1.00 50.00 .
ATOM   369  ND2 ND2 . ASN ASN . . 54  54  . 13.913  35.749 8.208   1.00 50.00 .
ATOM   370  N   N   . SER SER . . 55  55  . 11.456  35.029 3.173   1.00 50.00 .
ATOM   371  CA  CA  . SER SER . . 55  55  . 10.308  34.917 2.251   1.00 50.00 .
ATOM   372  C   C   . SER SER . . 55  55  . 10.687  34.221 0.932   1.00 50.00 .
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ATOM   373  O   O   . SER SER . . 55  55  . 9.943   34.269 -0.048  1.00 50.00 .
ATOM   374  CB  CB  . SER SER . . 55  55  . 9.731   36.308 1.948   1.00 50.00 .
ATOM   375  OG  OG  . SER SER . . 55  55  . 9.327   36.967 3.152   1.00 50.00 .
ATOM   376  N   N   . TYR TYR . . 56  56  . 11.871  33.604 0.908   1.00 50.00 .
ATOM   377  CA  CA  . TYR TYR . . 56  56  . 12.404  32.986 -0.317  1.00 50.00 .
ATOM   378  C   C   . TYR TYR . . 56  56  . 12.255  31.473 -0.212  1.00 50.00 .
ATOM   379  O   O   . TYR TYR . . 56  56  . 12.575  30.851 0.804   1.00 50.00 .
ATOM   380  CB  CB  . TYR TYR . . 56  56  . 13.877  33.354 -0.547  1.00 50.00 .
ATOM   381  CG  CG  . TYR TYR . . 56  56  . 14.147  34.828 -0.880  1.00 50.00 .
ATOM   382  CD1 CD1 . TYR TYR . . 56  56  . 13.377  35.854 -0.345  1.00 50.00 .
ATOM   383  CD2 CD2 . TYR TYR . . 56  56  . 15.207  35.155 -1.718  1.00 50.00 .
ATOM   384  CE1 CE1 . TYR TYR . . 56  56  . 13.638  37.174 -0.622  1.00 50.00 .
ATOM   385  CE2 CE2 . TYR TYR . . 56  56  . 15.473  36.486 -2.006  1.00 50.00 .
ATOM   386  CZ  CZ  . TYR TYR . . 56  56  . 14.689  37.488 -1.458  1.00 50.00 .
ATOM   387  OH  OH  . TYR TYR . . 56  56  . 14.943  38.775 -1.763  1.00 50.00 .
ATOM   388  N   N   . TRP TRP . . 57  57  . 11.713  30.941 -1.294  1.00 50.00 .
ATOM   389  CA  CA  . TRP TRP . . 57  57  . 11.451  29.502 -1.425  1.00 50.00 .
ATOM   390  C   C   . TRP TRP . . 57  57  . 12.102  29.006 -2.716  1.00 50.00 .
ATOM   391  O   O   . TRP TRP . . 57  57  . 12.174  29.736 -3.707  1.00 50.00 .
ATOM   392  CB  CB  . TRP TRP . . 57  57  . 9.939   29.249 -1.474  1.00 50.00 .
ATOM   393  CG  CG  . TRP TRP . . 57  57  . 9.211   29.784 -0.236  1.00 50.00 .
ATOM   394  CD1 CD1 . TRP TRP . . 57  57  . 8.822   31.044 -0.046  1.00 50.00 .
ATOM   395  CD2 CD2 . TRP TRP . . 57  57  . 8.852   29.065 0.894   1.00 50.00 .
ATOM   396  NE1 NE1 . TRP TRP . . 57  57  . 8.231   31.156 1.143   1.00 50.00 .
ATOM   397  CE2 CE2 . TRP TRP . . 57  57  . 8.227   29.967 1.743   1.00 50.00 .
ATOM   398  CE3 CE3 . TRP TRP . . 57  57  . 8.952   27.723 1.239   1.00 50.00 .
ATOM   399  CZ2 CZ2 . TRP TRP . . 57  57  . 7.692   29.523 2.947   1.00 50.00 .
ATOM   400  CZ3 CZ3 . TRP TRP . . 57  57  . 8.405   27.277 2.436   1.00 50.00 .
ATOM   401  CH2 CH2 . TRP TRP . . 57  57  . 7.774   28.178 3.289   1.00 50.00 .
ATOM   402  N   N   . ARG ARG . . 58  58  . 12.596  27.778 -2.657  1.00 50.00 .
ATOM   403  CA  CA  . ARG ARG . . 58  58  . 13.191  27.125 -3.829  1.00 50.00 .
ATOM   404  C   C   . ARG ARG . . 58  58  . 12.311  25.954 -4.272  1.00 50.00 .
ATOM   405  O   O   . ARG ARG . . 58  58  . 11.918  25.083 -3.494  1.00 50.00 .
ATOM   406  CB  CB  . ARG ARG . . 58  58  . 14.614  26.675 -3.497  1.00 50.00 .
ATOM   407  CG  CG  . ARG ARG . . 58  58  . 15.310  25.964 -4.660  1.00 50.00 .
ATOM   408  CD  CD  . ARG ARG . . 58  58  . 16.678  25.430 -4.228  1.00 50.00 .
ATOM   409  NE  NE  . ARG ARG . . 58  58  . 17.577  26.536 -3.846  1.00 50.00 .
ATOM   410  CZ  CZ  . ARG ARG . . 58  58  . 18.301  27.291 -4.677  1.00 50.00 .
ATOM   411  NH1 NH1 . ARG ARG . . 58  58  . 18.260  27.106 -5.985  1.00 50.00 .
ATOM   412  NH2 NH2 . ARG ARG . . 58  58  . 19.030  28.295 -4.215  1.00 50.00 .
ATOM   413  N   N   . ILE ILE . . 59  59  . 11.974  26.014 -5.550  1.00 50.00 .
ATOM   414  CA  CA  . ILE ILE . . 59  59  . 11.165  24.973 -6.208  1.00 50.00 .
ATOM   415  C   C   . ILE ILE . . 59  59  . 12.006  23.704 -6.370  1.00 50.00 .
ATOM   416  O   O   . ILE ILE . . 59  59  . 13.046  23.710 -7.014  1.00 50.00 .
ATOM   417  CB  CB  . ILE ILE . . 59  59  . 10.675  25.514 -7.557  1.00 50.00 .
ATOM   418  CG1 CG1 . ILE ILE . . 59  59  . 9.784   26.725 -7.266  1.00 50.00 .
ATOM   419  CG2 CG2 . ILE ILE . . 59  59  . 9.969   24.454 -8.417  1.00 50.00 .
ATOM   420  CD1 CD1 . ILE ILE . . 59  59  . 9.309   27.433 -8.518  1.00 50.00 .
ATOM   421  N   N   . ARG ARG . . 60  60  . 11.461  22.607 -5.884  1.00 50.00 .
ATOM   422  CA  CA  . ARG ARG . . 60  60  . 12.085  21.284 -6.039  1.00 50.00 .
ATOM   423  C   C   . ARG ARG . . 60  60  . 11.138  20.338 -6.760  1.00 50.00 .
ATOM   424  O   O   . ARG ARG . . 60  60  . 9.922   20.448 -6.606  1.00 50.00 .
ATOM   425  CB  CB  . ARG ARG . . 60  60  . 12.385  20.696 -4.670  1.00 50.00 .
ATOM   426  CG  CG  . ARG ARG . . 60  60  . 13.475  21.491 -3.966  1.00 50.00 .
ATOM   427  CD  CD  . ARG ARG . . 60  60  . 13.681  20.847 -2.605  1.00 50.00 .
ATOM   428  NE  NE  . ARG ARG . . 60  60  . 14.294  19.516 -2.670  1.00 50.00 .
ATOM   429  CZ  CZ  . ARG ARG . . 60  60  . 15.596  19.267 -2.810  1.00 50.00 .
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ATOM   430  NH1 NH1 . ARG ARG . . 60  60  . 16.498  20.239 -2.885  1.00 50.00 .
ATOM   431  NH2 NH2 . ARG ARG . . 60  60  . 16.001  18.008 -2.872  1.00 50.00 .
ATOM   432  N   N   . GLY GLY . . 61  61  . 11.706  19.324 -7.421  1.00 50.00 .
ATOM   433  CA  CA  . GLY GLY . . 61  61  . 10.887  18.281 -8.068  1.00 50.00 .
ATOM   434  C   C   . GLY GLY . . 61  61  . 10.036  17.565 -7.006  1.00 50.00 .
ATOM   435  O   O   . GLY GLY . . 61  61  . 10.364  17.602 -5.818  1.00 50.00 .
ATOM   436  N   N   . GLY GLY . . 62  62  . 8.918   16.972 -7.444  1.00 50.00 .
ATOM   437  CA  CA  . GLY GLY . . 62  62  . 8.080   16.133 -6.558  1.00 50.00 .
ATOM   438  C   C   . GLY GLY . . 62  62  . 8.964   15.105 -5.829  1.00 50.00 .
ATOM   439  O   O   . GLY GLY . . 62  62  . 10.024  14.733 -6.331  1.00 50.00 .
ATOM   440  N   N   . SER SER . . 63  63  . 8.504   14.639 -4.670  1.00 50.00 .
ATOM   441  CA  CA  . SER SER . . 63  63  . 9.281   13.690 -3.829  1.00 50.00 .
ATOM   442  C   C   . SER SER . . 63  63  . 9.829   12.487 -4.608  1.00 50.00 .
ATOM   443  O   O   . SER SER . . 63  63  . 11.006  12.160 -4.504  1.00 50.00 .
ATOM   444  CB  CB  . SER SER . . 63  63  . 8.428   13.131 -2.692  1.00 50.00 .
ATOM   445  OG  OG  . SER SER . . 63  63  . 9.113   12.205 -1.860  1.00 50.00 .
ATOM   446  N   N   . GLU GLU . . 64  64  . 8.979   11.921 -5.463  1.00 50.00 .
ATOM   447  CA  CA  . GLU GLU . . 64  64  . 9.344   10.764 -6.300  1.00 50.00 .
ATOM   448  C   C   . GLU GLU . . 64  64  . 9.665   11.135 -7.758  1.00 50.00 .
ATOM   449  O   O   . GLU GLU . . 64  64  . 9.519   10.324 -8.672  1.00 50.00 .
ATOM   450  CB  CB  . GLU GLU . . 64  64  . 8.206   9.738  -6.232  1.00 50.00 .
ATOM   451  CG  CG  . GLU GLU . . 64  64  . 8.103   9.075  -4.853  1.00 50.00 .
ATOM   452  CD  CD  . GLU GLU . . 64  64  . 9.307   8.207  -4.454  1.00 50.00 .
ATOM   453  OE1 OE1 . GLU GLU . . 64  64  . 10.252  8.054  -5.260  1.00 50.00 .
ATOM   454  OE2 OE2 . GLU GLU . . 64  64  . 9.221   7.657  -3.335  1.00 50.00 .
ATOM   455  N   N   . GLY GLY . . 65  65  . 10.077  12.393 -7.943  1.00 50.00 .
ATOM   456  CA  CA  . GLY GLY . . 65  65  . 10.452  12.931 -9.259  1.00 50.00 .
ATOM   457  C   C   . GLY GLY . . 65  65  . 11.892  13.451 -9.233  1.00 50.00 .
ATOM   458  O   O   . GLY GLY . . 65  65  . 12.557  13.479 -8.198  1.00 50.00 .
ATOM   459  N   N   . GLY GLY . . 66  66  . 12.321  13.889 -10.415 1.00 50.00 .
ATOM   460  CA  CA  . GLY GLY . . 66  66  . 13.648  14.495 -10.612 1.00 50.00 .
ATOM   461  C   C   . GLY GLY . . 66  66  . 13.561  15.508 -11.754 1.00 50.00 .
ATOM   462  O   O   . GLY GLY . . 66  66  . 12.724  15.376 -12.647 1.00 50.00 .
ATOM   463  N   N   . CYS CYS . . 67  67  . 14.369  16.551 -11.635 1.00 50.00 .
ATOM   464  CA  CA  . CYS CYS . . 67  67  . 14.416  17.628 -12.636 1.00 50.00 .
ATOM   465  C   C   . CYS CYS . . 67  67  . 15.846  18.181 -12.751 1.00 50.00 .
ATOM   466  O   O   . CYS CYS . . 67  67  . 16.621  18.070 -11.796 1.00 50.00 .
ATOM   467  CB  CB  . CYS CYS . . 67  67  . 13.467  18.760 -12.225 1.00 50.00 .
ATOM   468  SG  SG  . CYS CYS . . 67  67  . 14.093  19.714 -10.792 1.00 50.00 .
ATOM   469  N   N   . PRO PRO . . 68  68  . 16.181  18.724 -13.920 1.00 50.00 .
ATOM   470  CA  CA  . PRO PRO . . 68  68  . 17.378  19.566 -14.049 1.00 50.00 .
ATOM   471  C   C   . PRO PRO . . 68  68  . 17.010  21.044 -13.878 1.00 50.00 .
ATOM   472  O   O   . PRO PRO . . 68  68  . 15.973  21.536 -14.308 1.00 50.00 .
ATOM   473  CB  CB  . PRO PRO . . 68  68  . 17.926  19.238 -15.440 1.00 50.00 .
ATOM   474  CG  CG  . PRO PRO . . 68  68  . 16.704  18.813 -16.252 1.00 50.00 .
ATOM   475  CD  CD  . PRO PRO . . 68  68  . 15.729  18.220 -15.229 1.00 50.00 .
ATOM   476  N   N   . ARG ARG . . 69  69  . 17.853  21.748 -13.143 1.00 50.00 .
ATOM   477  CA  CA  . ARG ARG . . 69  69  . 17.799  23.220 -12.996 1.00 50.00 .
ATOM   478  C   C   . ARG ARG . . 69  69  . 17.640  23.913 -14.361 1.00 50.00 .
ATOM   479  O   O   . ARG ARG . . 69  69  . 18.279  23.533 -15.343 1.00 50.00 .
ATOM   480  CB  CB  . ARG ARG . . 69  69  . 19.068  23.757 -12.305 1.00 50.00 .
ATOM   481  CG  CG  . ARG ARG . . 69  69  . 20.250  22.776 -12.182 1.00 50.00 .
ATOM   482  CD  CD  . ARG ARG . . 69  69  . 20.762  22.299 -13.545 1.00 50.00 .
ATOM   483  NE  NE  . ARG ARG . . 69  69  . 21.464  21.013 -13.412 1.00 50.00 .
ATOM   484  CZ  CZ  . ARG ARG . . 69  69  . 21.971  20.309 -14.426 1.00 50.00 .
ATOM   485  NH1 NH1 . ARG ARG . . 69  69  . 21.849  20.724 -15.682 1.00 50.00 .
ATOM   486  NH2 NH2 . ARG ARG . . 69  69  . 22.618  19.177 -14.191 1.00 50.00 .
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ATOM   487  N   N   . GLY GLY . . 70  70  . 16.653  24.813 -14.414 1.00 50.00 .
ATOM   488  CA  CA  . GLY GLY . . 70  70  . 16.290  25.482 -15.680 1.00 50.00 .
ATOM   489  C   C   . GLY GLY . . 70  70  . 15.086  24.821 -16.371 1.00 50.00 .
ATOM   490  O   O   . GLY GLY . . 70  70  . 14.468  25.420 -17.248 1.00 50.00 .
ATOM   491  N   N   . SER SER . . 71  71  . 14.750  23.600 -15.955 1.00 50.00 .
ATOM   492  CA  CA  . SER SER . . 71  71  . 13.591  22.880 -16.509 1.00 50.00 .
ATOM   493  C   C   . SER SER . . 71  71  . 12.291  23.616 -16.215 1.00 50.00 .
ATOM   494  O   O   . SER SER . . 71  71  . 12.059  23.981 -15.062 1.00 50.00 .
ATOM   495  CB  CB  . SER SER . . 71  71  . 13.436  21.477 -15.943 1.00 50.00 .
ATOM   496  OG  OG  . SER SER . . 71  71  . 14.491  20.718 -16.507 1.00 50.00 .
ATOM   497  N   N   . PRO PRO . . 72  72  . 11.477  23.821 -17.251 1.00 50.00 .
ATOM   498  CA  CA  . PRO PRO . . 72  72  . 10.151  24.425 -17.090 1.00 50.00 .
ATOM   499  C   C   . PRO PRO . . 72  72  . 9.266   23.446 -16.315 1.00 50.00 .
ATOM   500  O   O   . PRO PRO . . 72  72  . 9.394   22.227 -16.416 1.00 50.00 .
ATOM   501  CB  CB  . PRO PRO . . 72  72  . 9.656   24.622 -18.521 1.00 50.00 .
ATOM   502  CG  CG  . PRO PRO . . 72  72  . 10.312  23.476 -19.290 1.00 50.00 .
ATOM   503  CD  CD  . PRO PRO . . 72  72  . 11.694  23.380 -18.646 1.00 50.00 .
ATOM   504  N   N   . VAL VAL . . 73  73  . 8.409   24.025 -15.502 1.00 50.00 .
ATOM   505  CA  CA  . VAL VAL . . 73  73  . 7.439   23.248 -14.711 1.00 50.00 .
ATOM   506  C   C   . VAL VAL . . 73  73  . 6.080   23.335 -15.421 1.00 50.00 .
ATOM   507  O   O   . VAL VAL . . 73  73  . 5.587   24.410 -15.758 1.00 50.00 .
ATOM   508  CB  CB  . VAL VAL . . 73  73  . 7.381   23.798 -13.274 1.00 50.00 .
ATOM   509  CG1 CG1 . VAL VAL . . 73  73  . 8.697   23.627 -12.510 1.00 50.00 .
ATOM   510  CG2 CG2 . VAL VAL . . 73  73  . 7.151   25.298 -13.315 1.00 50.00 .
ATOM   511  N   N   . ARG ARG . . 74  74  . 5.529   22.183 -15.748 1.00 50.00 .
ATOM   512  CA  CA  . ARG ARG . . 74  74  . 4.228   22.094 -16.428 1.00 50.00 .
ATOM   513  C   C   . ARG ARG . . 74  74  . 3.108   22.144 -15.387 1.00 50.00 .
ATOM   514  O   O   . ARG ARG . . 74  74  . 3.217   21.613 -14.280 1.00 50.00 .
ATOM   515  CB  CB  . ARG ARG . . 74  74  . 4.123   20.798 -17.239 1.00 50.00 .
ATOM   516  CG  CG  . ARG ARG . . 74  74  . 5.189   20.702 -18.331 1.00 50.00 .
ATOM   517  CD  CD  . ARG ARG . . 74  74  . 4.993   19.420 -19.141 1.00 50.00 .
ATOM   518  NE  NE  . ARG ARG . . 74  74  . 5.989   19.343 -20.226 1.00 50.00 .
ATOM   519  CZ  CZ  . ARG ARG . . 74  74  . 7.257   18.932 -20.119 1.00 50.00 .
ATOM   520  NH1 NH1 . ARG ARG . . 74  74  . 7.759   18.533 -18.958 1.00 50.00 .
ATOM   521  NH2 NH2 . ARG ARG . . 74  74  . 8.035   18.895 -21.192 1.00 50.00 .
ATOM   522  N   N   . CYS CYS . . 75  75  . 2.036   22.826 -15.765 1.00 50.00 .
ATOM   523  CA  CA  . CYS CYS . . 75  75  . 0.819   22.906 -14.936 1.00 50.00 .
ATOM   524  C   C   . CYS CYS . . 75  75  . 0.235   21.499 -14.746 1.00 50.00 .
ATOM   525  O   O   . CYS CYS . . 75  75  . 0.262   20.683 -15.669 1.00 50.00 .
ATOM   526  CB  CB  . CYS CYS . . 75  75  . -0.215  23.785 -15.634 1.00 50.00 .
ATOM   527  SG  SG  . CYS CYS . . 75  75  . 0.377   25.484 -15.974 1.00 50.00 .
ATOM   528  N   N   . GLY GLY . . 76  76  . -0.094  21.199 -13.487 1.00 50.00 .
ATOM   529  CA  CA  . GLY GLY . . 76  76  . -0.626  19.874 -13.117 1.00 50.00 .
ATOM   530  C   C   . GLY GLY . . 76  76  . 0.432   19.015 -12.411 1.00 50.00 .
ATOM   531  O   O   . GLY GLY . . 76  76  . 0.099   18.086 -11.678 1.00 50.00 .
ATOM   532  N   N   . GLN GLN . . 77  77  . 1.701   19.315 -12.685 1.00 50.00 .
ATOM   533  CA  CA  . GLN GLN . . 77  77  . 2.832   18.625 -12.042 1.00 50.00 .
ATOM   534  C   C   . GLN GLN . . 77  77  . 2.942   19.029 -10.567 1.00 50.00 .
ATOM   535  O   O   . GLN GLN . . 77  77  . 2.569   20.136 -10.167 1.00 50.00 .
ATOM   536  CB  CB  . GLN GLN . . 77  77  . 4.143   18.934 -12.768 1.00 50.00 .
ATOM   537  CG  CG  . GLN GLN . . 77  77  . 4.145   18.381 -14.196 1.00 50.00 .
ATOM   538  CD  CD  . GLN GLN . . 77  77  . 5.496   18.619 -14.869 1.00 50.00 .
ATOM   539  OE1 OE1 . GLN GLN . . 77  77  . 6.018   19.718 -14.968 1.00 50.00 .
ATOM   540  NE2 NE2 . GLN GLN . . 77  77  . 6.151   17.569 -15.284 1.00 50.00 .
ATOM   541  N   N   . ALA ALA . . 78  78  . 3.387   18.060 -9.777  1.00 50.00 .
ATOM   542  CA  CA  . ALA ALA . . 78  78  . 3.647   18.262 -8.344  1.00 50.00 .
ATOM   543  C   C   . ALA ALA . . 78  78  . 5.079   18.773 -8.135  1.00 50.00 .
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ATOM   544  O   O   . ALA ALA . . 78  78  . 6.001   18.443 -8.880  1.00 50.00 .
ATOM   545  CB  CB  . ALA ALA . . 78  78  . 3.433   16.952 -7.583  1.00 50.00 .
ATOM   546  N   N   . VAL VAL . . 79  79  . 5.207   19.604 -7.115  1.00 50.00 .
ATOM   547  CA  CA  . VAL VAL . . 79  79  . 6.472   20.228 -6.701  1.00 50.00 .
ATOM   548  C   C   . VAL VAL . . 79  79  . 6.582   20.204 -5.164  1.00 50.00 .
ATOM   549  O   O   . VAL VAL . . 79  79  . 5.610   19.935 -4.459  1.00 50.00 .
ATOM   550  CB  CB  . VAL VAL . . 79  79  . 6.497   21.655 -7.292  1.00 50.00 .
ATOM   551  CG1 CG1 . VAL VAL . . 79  79  . 5.448   22.598 -6.700  1.00 50.00 .
ATOM   552  CG2 CG2 . VAL VAL . . 79  79  . 7.835   22.341 -7.109  1.00 50.00 .
ATOM   553  N   N   . ARG ARG . . 80  80  . 7.789   20.483 -4.700  1.00 50.00 .
ATOM   554  CA  CA  . ARG ARG . . 80  80  . 8.071   20.793 -3.291  1.00 50.00 .
ATOM   555  C   C   . ARG ARG . . 80  80  . 8.680   22.196 -3.200  1.00 50.00 .
ATOM   556  O   O   . ARG ARG . . 80  80  . 9.340   22.666 -4.126  1.00 50.00 .
ATOM   557  CB  CB  . ARG ARG . . 80  80  . 9.040   19.765 -2.711  1.00 50.00 .
ATOM   558  CG  CG  . ARG ARG . . 80  80  . 8.370   18.395 -2.586  1.00 50.00 .
ATOM   559  CD  CD  . ARG ARG . . 80  80  . 9.348   17.362 -2.033  1.00 50.00 .
ATOM   560  NE  NE  . ARG ARG . . 80  80  . 10.411  17.106 -3.018  1.00 50.00 .
ATOM   561  CZ  CZ  . ARG ARG . . 80  80  . 11.613  16.590 -2.765  1.00 50.00 .
ATOM   562  NH1 NH1 . ARG ARG . . 80  80  . 11.987  16.270 -1.534  1.00 50.00 .
ATOM   563  NH2 NH2 . ARG ARG . . 80  80  . 12.459  16.393 -3.765  1.00 50.00 .
ATOM   564  N   N   . LEU LEU . . 81  81  . 8.408   22.855 -2.081  1.00 50.00 .
ATOM   565  CA  CA  . LEU LEU . . 81  81  . 8.899   24.221 -1.831  1.00 50.00 .
ATOM   566  C   C   . LEU LEU . . 81  81  . 9.782   24.284 -0.589  1.00 50.00 .
ATOM   567  O   O   . LEU LEU . . 81  81  . 9.319   24.001 0.517   1.00 50.00 .
ATOM   568  CB  CB  . LEU LEU . . 81  81  . 7.737   25.210 -1.688  1.00 50.00 .
ATOM   569  CG  CG  . LEU LEU . . 81  81  . 6.990   25.449 -3.004  1.00 50.00 .
ATOM   570  CD1 CD1 . LEU LEU . . 81  81  . 5.779   26.341 -2.738  1.00 50.00 .
ATOM   571  CD2 CD2 . LEU LEU . . 81  81  . 7.895   26.108 -4.049  1.00 50.00 .
ATOM   572  N   N   . THR THR . . 82  82  . 11.077  24.449 -0.823  1.00 50.00 .
ATOM   573  CA  CA  . THR THR . . 82  82  . 12.042  24.606 0.282   1.00 50.00 .
ATOM   574  C   C   . THR THR . . 82  82  . 12.106  26.046 0.766   1.00 50.00 .
ATOM   575  O   O   . THR THR . . 82  82  . 12.287  26.978 -0.011  1.00 50.00 .
ATOM   576  CB  CB  . THR THR . . 82  82  . 13.483  24.237 -0.072  1.00 50.00 .
ATOM   577  OG1 OG1 . THR THR . . 82  82  . 13.833  24.772 -1.341  1.00 50.00 .
ATOM   578  CG2 CG2 . THR THR . . 82  82  . 13.739  22.747 0.002   1.00 50.00 .
ATOM   579  N   N   . HIS HIS . . 83  83  . 12.041  26.171 2.079   1.00 50.00 .
ATOM   580  CA  CA  . HIS HIS . . 83  83  . 12.283  27.448 2.757   1.00 50.00 .
ATOM   581  C   C   . HIS HIS . . 83  83  . 13.800  27.584 2.923   1.00 50.00 .
ATOM   582  O   O   . HIS HIS . . 83  83  . 14.403  26.894 3.744   1.00 50.00 .
ATOM   583  CB  CB  . HIS HIS . . 83  83  . 11.556  27.447 4.104   1.00 50.00 .
ATOM   584  CG  CG  . HIS HIS . . 83  83  . 11.721  28.786 4.823   1.00 50.00 .
ATOM   585  ND1 ND1 . HIS HIS . . 83  83  . 12.366  28.966 5.970   1.00 50.00 .
ATOM   586  CD2 CD2 . HIS HIS . . 83  83  . 11.293  29.970 4.400   1.00 50.00 .
ATOM   587  CE1 CE1 . HIS HIS . . 83  83  . 12.338  30.256 6.277   1.00 50.00 .
ATOM   588  NE2 NE2 . HIS HIS . . 83  83  . 11.680  30.865 5.300   1.00 50.00 .
ATOM   589  N   N   . VAL VAL . . 84  84  . 14.385  28.489 2.147   1.00 50.00 .
ATOM   590  CA  CA  . VAL VAL . . 84  84  . 15.855  28.649 2.051   1.00 50.00 .
ATOM   591  C   C   . VAL VAL . . 84  84  . 16.580  28.738 3.409   1.00 50.00 .
ATOM   592  O   O   . VAL VAL . . 84  84  . 17.528  27.997 3.656   1.00 50.00 .
ATOM   593  CB  CB  . VAL VAL . . 84  84  . 16.204  29.851 1.144   1.00 50.00 .
ATOM   594  CG1 CG1 . VAL VAL . . 84  84  . 15.657  31.187 1.665   1.00 50.00 .
ATOM   595  CG2 CG2 . VAL VAL . . 84  84  . 17.713  29.962 0.896   1.00 50.00 .
ATOM   596  N   N   . LEU LEU . . 85  85  . 16.072  29.602 4.284   1.00 50.00 .
ATOM   597  CA  CA  . LEU LEU . . 85  85  . 16.779  29.940 5.531   1.00 50.00 .
ATOM   598  C   C   . LEU LEU . . 85  85  . 16.781  28.793 6.554   1.00 50.00 .
ATOM   599  O   O   . LEU LEU . . 85  85  . 17.701  28.678 7.359   1.00 50.00 .
ATOM   600  CB  CB  . LEU LEU . . 85  85  . 16.253  31.264 6.107   1.00 50.00 .
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ATOM   601  CG  CG  . LEU LEU . . 85  85  . 17.081  31.732 7.314   1.00 50.00 .
ATOM   602  CD1 CD1 . LEU LEU . . 85  85  . 17.208  33.255 7.345   1.00 50.00 .
ATOM   603  CD2 CD2 . LEU LEU . . 85  85  . 16.455  31.260 8.629   1.00 50.00 .
ATOM   604  N   N   . THR THR . . 86  86  . 15.741  27.967 6.524   1.00 50.00 .
ATOM   605  CA  CA  . THR THR . . 86  86  . 15.599  26.878 7.512   1.00 50.00 .
ATOM   606  C   C   . THR THR . . 86  86  . 15.786  25.471 6.919   1.00 50.00 .
ATOM   607  O   O   . THR THR . . 86  86  . 15.779  24.488 7.656   1.00 50.00 .
ATOM   608  CB  CB  . THR THR . . 86  86  . 14.262  26.971 8.267   1.00 50.00 .
ATOM   609  OG1 OG1 . THR THR . . 86  86  . 13.163  26.784 7.369   1.00 50.00 .
ATOM   610  CG2 CG2 . THR THR . . 86  86  . 14.120  28.302 9.017   1.00 50.00 .
ATOM   611  N   N   . GLY GLY . . 87  87  . 15.849  25.404 5.581   1.00 50.00 .
ATOM   612  CA  CA  . GLY GLY . . 87  87  . 15.968  24.148 4.811   1.00 50.00 .
ATOM   613  C   C   . GLY GLY . . 87  87  . 14.783  23.182 4.988   1.00 50.00 .
ATOM   614  O   O   . GLY GLY . . 87  87  . 14.881  22.011 4.627   1.00 50.00 .
ATOM   615  N   N   . LYS LYS . . 88  88  . 13.673  23.710 5.499   1.00 50.00 .
ATOM   616  CA  CA  . LYS LYS . . 88  88  . 12.409  22.968 5.651   1.00 50.00 .
ATOM   617  C   C   . LYS LYS . . 88  88  . 11.636  23.007 4.323   1.00 50.00 .
ATOM   618  O   O   . LYS LYS . . 88  88  . 11.988  23.733 3.392   1.00 50.00 .
ATOM   619  CB  CB  . LYS LYS . . 88  88  . 11.581  23.608 6.775   1.00 50.00 .
ATOM   620  CG  CG  . LYS LYS . . 88  88  . 12.243  23.553 8.159   1.00 50.00 .
ATOM   621  CD  CD  . LYS LYS . . 88  88  . 12.208  22.144 8.746   1.00 50.00 .
ATOM   622  CE  CE  . LYS LYS . . 88  88  . 13.014  21.956 10.035  1.00 50.00 .
ATOM   623  NZ  NZ  . LYS LYS . . 88  88  . 14.439  22.297 9.890   1.00 50.00 .
ATOM   624  N   N   . ASN ASN . . 89  89  . 10.563  22.233 4.263   1.00 50.00 .
ATOM   625  CA  CA  . ASN ASN . . 89  89  . 9.670   22.194 3.094   1.00 50.00 .
ATOM   626  C   C   . ASN ASN . . 89  89  . 8.244   22.558 3.491   1.00 50.00 .
ATOM   627  O   O   . ASN ASN . . 89  89  . 7.788   22.190 4.574   1.00 50.00 .
ATOM   628  CB  CB  . ASN ASN . . 89  89  . 9.671   20.798 2.464   1.00 50.00 .
ATOM   629  CG  CG  . ASN ASN . . 89  89  . 11.019  20.489 1.818   1.00 50.00 .
ATOM   630  OD1 OD1 . ASN ASN . . 89  89  . 11.755  19.591 2.198   1.00 50.00 .
ATOM   631  ND2 ND2 . ASN ASN . . 89  89  . 11.318  21.233 0.783   1.00 50.00 .
ATOM   632  N   N   . LEU LEU . . 90  90  . 7.565   23.243 2.577   1.00 50.00 .
ATOM   633  CA  CA  . LEU LEU . . 90  90  . 6.132   23.550 2.715   1.00 50.00 .
ATOM   634  C   C   . LEU LEU . . 90  90  . 5.359   22.230 2.844   1.00 50.00 .
ATOM   635  O   O   . LEU LEU . . 90  90  . 5.403   21.378 1.959   1.00 50.00 .
ATOM   636  CB  CB  . LEU LEU . . 90  90  . 5.642   24.333 1.494   1.00 50.00 .
ATOM   637  CG  CG  . LEU LEU . . 90  90  . 4.236   24.904 1.709   1.00 50.00 .
ATOM   638  CD1 CD1 . LEU LEU . . 90  90  . 4.242   25.983 2.795   1.00 50.00 .
ATOM   639  CD2 CD2 . LEU LEU . . 90  90  . 3.692   25.470 0.398   1.00 50.00 .
ATOM   640  N   N   . HIS HIS . . 91  91  . 4.814   22.033 4.033   1.00 50.00 .
ATOM   641  CA  CA  . HIS HIS . . 91  91  . 4.214   20.757 4.428   1.00 50.00 .
ATOM   642  C   C   . HIS HIS . . 91  91  . 2.773   20.943 4.890   1.00 50.00 .
ATOM   643  O   O   . HIS HIS . . 91  91  . 2.357   22.004 5.366   1.00 50.00 .
ATOM   644  CB  CB  . HIS HIS . . 91  91  . 5.026   20.179 5.579   1.00 50.00 .
ATOM   645  CG  CG  . HIS HIS . . 91  91  . 4.695   18.704 5.803   1.00 50.00 .
ATOM   646  ND1 ND1 . HIS HIS . . 91  91  . 5.397   17.708 5.293   1.00 50.00 .
ATOM   647  CD2 CD2 . HIS HIS . . 91  91  . 3.863   18.204 6.709   1.00 50.00 .
ATOM   648  CE1 CE1 . HIS HIS . . 91  91  . 5.071   16.601 5.944   1.00 50.00 .
ATOM   649  NE2 NE2 . HIS HIS . . 91  91  . 4.092   16.898 6.788   1.00 50.00 .
ATOM   650  N   N   . THR THR . . 92  92  . 2.030   19.861 4.733   1.00 50.00 .
ATOM   651  CA  CA  . THR THR . . 92  92  . 0.627   19.806 5.159   1.00 50.00 .
ATOM   652  C   C   . THR THR . . 92  92  . 0.256   18.370 5.546   1.00 50.00 .
ATOM   653  O   O   . THR THR . . 92  92  . 0.910   17.414 5.145   1.00 50.00 .
ATOM   654  CB  CB  . THR THR . . 92  92  . -0.249  20.329 4.024   1.00 50.00 .
ATOM   655  OG1 OG1 . THR THR . . 92  92  . 0.176   21.613 3.621   1.00 50.00 .
ATOM   656  CG2 CG2 . THR THR . . 92  92  . -1.608  20.681 4.558   1.00 50.00 .
ATOM   657  N   N   . HIS HIS . . 93  93  . -0.831  18.250 6.303   1.00 50.00 .
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ATOM   658  CA  CA  . HIS HIS . . 93  93  . -1.246  16.991 6.958   1.00 50.00 .
ATOM   659  C   C   . HIS HIS . . 93  93  . -2.599  17.147 7.673   1.00 50.00 .
ATOM   660  O   O   . HIS HIS . . 93  93  . -3.162  18.237 7.706   1.00 50.00 .
ATOM   661  CB  CB  . HIS HIS . . 93  93  . -0.161  16.547 7.963   1.00 50.00 .
ATOM   662  CG  CG  . HIS HIS . . 93  93  . 0.483   17.681 8.778   1.00 50.00 .
ATOM   663  ND1 ND1 . HIS HIS . . 93  93  . 1.655   17.597 9.390   1.00 50.00 .
ATOM   664  CD2 CD2 . HIS HIS . . 93  93  . 0.042   18.932 8.938   1.00 50.00 .
ATOM   665  CE1 CE1 . HIS HIS . . 93  93  . 1.946   18.793 9.886   1.00 50.00 .
ATOM   666  NE2 NE2 . HIS HIS . . 93  93  . 0.946   19.619 9.607   1.00 50.00 .
ATOM   667  N   N   . HIS HIS . . 94  94  . -2.993  16.107 8.400   1.00 50.00 .
ATOM   668  CA  CA  . HIS HIS . . 94  94  . -4.275  16.051 9.127   1.00 50.00 .
ATOM   669  C   C   . HIS HIS . . 94  94  . -4.243  16.658 10.540  1.00 50.00 .
ATOM   670  O   O   . HIS HIS . . 94  94  . -4.671  16.058 11.524  1.00 50.00 .
ATOM   671  CB  CB  . HIS HIS . . 94  94  . -4.782  14.607 9.139   1.00 50.00 .
ATOM   672  CG  CG  . HIS HIS . . 94  94  . -5.517  14.263 7.844   1.00 50.00 .
ATOM   673  ND1 ND1 . HIS HIS . . 94  94  . -6.432  15.013 7.235   1.00 50.00 .
ATOM   674  CD2 CD2 . HIS HIS . . 94  94  . -5.353  13.150 7.138   1.00 50.00 .
ATOM   675  CE1 CE1 . HIS HIS . . 94  94  . -6.836  14.360 6.150   1.00 50.00 .
ATOM   676  NE2 NE2 . HIS HIS . . 94  94  . -6.167  13.212 6.090   1.00 50.00 .
ATOM   677  N   N   . PHE PHE . . 95  95  . -3.720  17.876 10.606  1.00 50.00 .
ATOM   678  CA  CA  . PHE PHE . . 95  95  . -3.749  18.681 11.837  1.00 50.00 .
ATOM   679  C   C   . PHE PHE . . 95  95  . -4.610  19.933 11.619  1.00 50.00 .
ATOM   680  O   O   . PHE PHE . . 95  95  . -4.651  20.446 10.495  1.00 50.00 .
ATOM   681  CB  CB  . PHE PHE . . 95  95  . -2.337  19.078 12.281  1.00 50.00 .
ATOM   682  CG  CG  . PHE PHE . . 95  95  . -1.528  17.860 12.727  1.00 50.00 .
ATOM   683  CD1 CD1 . PHE PHE . . 95  95  . -1.517  17.435 14.047  1.00 50.00 .
ATOM   684  CD2 CD2 . PHE PHE . . 95  95  . -0.749  17.198 11.797  1.00 50.00 .
ATOM   685  CE1 CE1 . PHE PHE . . 95  95  . -0.709  16.364 14.414  1.00 50.00 .
ATOM   686  CE2 CE2 . PHE PHE . . 95  95  . 0.066   16.137 12.150  1.00 50.00 .
ATOM   687  CZ  CZ  . PHE PHE . . 95  95  . 0.085   15.718 13.472  1.00 50.00 .
ATOM   688  N   N   . PRO PRO . . 96  96  . -5.321  20.379 12.663  1.00 50.00 .
ATOM   689  CA  CA  . PRO PRO . . 96  96  . -6.113  21.617 12.615  1.00 50.00 .
ATOM   690  C   C   . PRO PRO . . 96  96  . -5.184  22.835 12.621  1.00 50.00 .
ATOM   691  O   O   . PRO PRO . . 96  96  . -4.134  22.857 13.263  1.00 50.00 .
ATOM   692  CB  CB  . PRO PRO . . 96  96  . -6.984  21.560 13.869  1.00 50.00 .
ATOM   693  CG  CG  . PRO PRO . . 96  96  . -6.119  20.789 14.866  1.00 50.00 .
ATOM   694  CD  CD  . PRO PRO . . 96  96  . -5.433  19.739 13.991  1.00 50.00 .
ATOM   695  N   N   . SER SER . . 97  97  . -5.573  23.817 11.827  1.00 50.00 .
ATOM   696  CA  CA  . SER SER . . 97  97  . -4.845  25.089 11.721  1.00 50.00 .
ATOM   697  C   C   . SER SER . . 97  97  . -4.816  25.815 13.078  1.00 50.00 .
ATOM   698  O   O   . SER SER . . 97  97  . -5.693  25.588 13.917  1.00 50.00 .
ATOM   699  CB  CB  . SER SER . . 97  97  . -5.537  25.903 10.639  1.00 50.00 .
ATOM   700  OG  OG  . SER SER . . 97  97  . -4.866  27.125 10.340  1.00 50.00 .
ATOM   701  N   N   . PRO PRO . . 98  98  . -3.761  26.605 13.329  1.00 50.00 .
ATOM   702  CA  CA  . PRO PRO . . 98  98  . -3.528  27.245 14.634  1.00 50.00 .
ATOM   703  C   C   . PRO PRO . . 98  98  . -4.664  28.162 15.100  1.00 50.00 .
ATOM   704  O   O   . PRO PRO . . 98  98  . -5.025  28.135 16.271  1.00 50.00 .
ATOM   705  CB  CB  . PRO PRO . . 98  98  . -2.203  27.992 14.465  1.00 50.00 .
ATOM   706  CG  CG  . PRO PRO . . 98  98  . -2.172  28.295 12.973  1.00 50.00 .
ATOM   707  CD  CD  . PRO PRO . . 98  98  . -2.705  27.004 12.376  1.00 50.00 .
ATOM   708  N   N   . LEU LEU . . 99  99  . -5.156  29.006 14.202  1.00 99.99 .
ATOM   709  CA  CA  . LEU LEU . . 99  99  . -6.201  29.984 14.545  1.00 99.99 .
ATOM   710  C   C   . LEU LEU . . 99  99  . -7.476  29.849 13.695  1.00 99.99 .
ATOM   711  O   O   . LEU LEU . . 99  99  . -8.562  30.236 14.112  1.00 99.99 .
ATOM   712  CB  CB  . LEU LEU . . 99  99  . -5.572  31.379 14.461  1.00 99.99 .
ATOM   713  CG  CG  . LEU LEU . . 99  99  . -4.791  31.870 15.687  1.00 99.99 .
ATOM   714  CD1 CD1 . LEU LEU . . 99  99  . -3.471  31.128 15.863  1.00 99.99 .
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ATOM   715  CD2 CD2 . LEU LEU . . 99  99  . -4.472  33.359 15.551  1.00 99.99 .
ATOM   716  N   N   . SER SER . . 100 100 . -7.336  29.264 12.511  1.00 99.99 .
ATOM   717  CA  CA  . SER SER . . 100 100 . -8.483  29.047 11.612  1.00 99.99 .
ATOM   718  C   C   . SER SER . . 100 100 . -8.482  27.643 11.068  1.00 99.99 .
ATOM   719  O   O   . SER SER . . 100 100 . -7.789  27.388 10.083  1.00 99.99 .
ATOM   720  CB  CB  . SER SER . . 100 100 . -8.467  30.078 10.483  1.00 99.99 .
ATOM   721  OG  OG  . SER SER . . 100 100 . -9.394  29.788 9.437   1.00 99.99 .
ATOM   722  N   N   . ASN ASN . . 101 101 . -9.007  26.736 11.884  1.00 99.99 .
ATOM   723  CA  CA  . ASN ASN . . 101 101 . -9.245  25.319 11.580  1.00 99.99 .
ATOM   724  C   C   . ASN ASN . . 101 101 . -10.317 25.066 10.511  1.00 99.99 .
ATOM   725  O   O   . ASN ASN . . 101 101 . -11.545 25.151 10.756  1.00 99.99 .
ATOM   726  CB  CB  . ASN ASN . . 101 101 . -9.615  24.625 12.907  1.00 99.99 .
ATOM   727  CG  CG  . ASN ASN . . 101 101 . -10.865 25.193 13.593  1.00 99.99 .
ATOM   728  OD1 OD1 . ASN ASN . . 101 101 . -11.742 25.880 13.057  1.00 99.99 .
ATOM   729  ND2 ND2 . ASN ASN . . 101 101 . -10.972 25.101 14.873  1.00 99.99 .
ATOM   730  N   N   . ASN ASN . . 102 102 . -10.116 24.746 9.254   1.00 50.00 .
ATOM   731  CA  CA  . ASN ASN . . 102 102 . -8.936  24.576 8.410   1.00 50.00 .
ATOM   732  C   C   . ASN ASN . . 102 102 . -7.783  23.692 8.881   1.00 50.00 .
ATOM   733  O   O   . ASN ASN . . 102 102 . -7.572  23.421 10.057  1.00 50.00 .
ATOM   734  CB  CB  . ASN ASN . . 102 102 . -8.404  25.904 7.875   1.00 50.00 .
ATOM   735  CG  CG  . ASN ASN . . 102 102 . -9.478  26.662 7.116   1.00 50.00 .
ATOM   736  OD1 OD1 . ASN ASN . . 102 102 . -9.832  27.807 7.352   1.00 50.00 .
ATOM   737  ND2 ND2 . ASN ASN . . 102 102 . -9.880  26.026 6.052   1.00 50.00 .
ATOM   738  N   N   . GLN GLN . . 103 103 . -7.136  23.183 7.853   1.00 50.00 .
ATOM   739  CA  CA  . GLN GLN . . 103 103 . -5.979  22.290 7.963   1.00 50.00 .
ATOM   740  C   C   . GLN GLN . . 103 103 . -4.703  23.124 8.138   1.00 50.00 .
ATOM   741  O   O   . GLN GLN . . 103 103 . -4.541  24.171 7.513   1.00 50.00 .
ATOM   742  CB  CB  . GLN GLN . . 103 103 . -5.960  21.482 6.667   1.00 50.00 .
ATOM   743  CG  CG  . GLN GLN . . 103 103 . -4.929  20.361 6.711   1.00 50.00 .
ATOM   744  CD  CD  . GLN GLN . . 103 103 . -5.073  19.431 5.514   1.00 50.00 .
ATOM   745  OE1 OE1 . GLN GLN . . 103 103 . -5.945  18.593 5.397   1.00 50.00 .
ATOM   746  NE2 NE2 . GLN GLN . . 103 103 . -4.193  19.599 4.556   1.00 50.00 .
ATOM   747  N   N   . GLU GLU . . 104 104 . -3.815  22.615 8.981   1.00 50.00 .
ATOM   748  CA  CA  . GLU GLU . . 104 104 . -2.520  23.255 9.268   1.00 50.00 .
ATOM   749  C   C   . GLU GLU . . 104 104 . -1.579  23.155 8.056   1.00 50.00 .
ATOM   750  O   O   . GLU GLU . . 104 104 . -1.514  22.125 7.385   1.00 50.00 .
ATOM   751  CB  CB  . GLU GLU . . 104 104 . -1.878  22.573 10.480  1.00 50.00 .
ATOM   752  CG  CG  . GLU GLU . . 104 104 . -0.633  23.329 10.952  1.00 50.00 .
ATOM   753  CD  CD  . GLU GLU . . 104 104 . 0.180   22.497 11.934  1.00 50.00 .
ATOM   754  OE1 OE1 . GLU GLU . . 104 104 . 1.027   21.736 11.425  1.00 50.00 .
ATOM   755  OE2 OE2 . GLU GLU . . 104 104 . -0.063  22.617 13.153  1.00 50.00 .
ATOM   756  N   N   . VAL VAL . . 105 105 . -0.804  24.219 7.877   1.00 50.00 .
ATOM   757  CA  CA  . VAL VAL . . 105 105 . 0.293   24.275 6.890   1.00 50.00 .
ATOM   758  C   C   . VAL VAL . . 105 105 . 1.578   24.645 7.642   1.00 50.00 .
ATOM   759  O   O   . VAL VAL . . 105 105 . 1.633   25.666 8.319   1.00 50.00 .
ATOM   760  CB  CB  . VAL VAL . . 105 105 . -0.018  25.299 5.784   1.00 50.00 .
ATOM   761  CG1 CG1 . VAL VAL . . 105 105 . 1.142   25.454 4.797   1.00 50.00 .
ATOM   762  CG2 CG2 . VAL VAL . . 105 105 . -1.230  24.848 4.973   1.00 50.00 .
ATOM   763  N   N   . SER SER . . 106 106 . 2.566   23.774 7.559   1.00 50.00 .
ATOM   764  CA  CA  . SER SER . . 106 106 . 3.845   23.956 8.270   1.00 50.00 .
ATOM   765  C   C   . SER SER . . 106 106 . 5.061   23.876 7.338   1.00 50.00 .
ATOM   766  O   O   . SER SER . . 106 106 . 4.929   23.690 6.131   1.00 50.00 .
ATOM   767  CB  CB  . SER SER . . 106 106 . 3.964   22.905 9.377   1.00 50.00 .
ATOM   768  OG  OG  . SER SER . . 106 106 . 3.857   21.602 8.801   1.00 50.00 .
ATOM   769  N   N   . ALA ALA . . 107 107 . 6.215   24.183 7.916   1.00 50.00 .
ATOM   770  CA  CA  . ALA ALA . . 107 107 . 7.530   24.035 7.271   1.00 50.00 .
ATOM   771  C   C   . ALA ALA . . 107 107 . 8.244   22.873 7.974   1.00 50.00 .
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ATOM   772  O   O   . ALA ALA . . 107 107 . 8.584   22.948 9.155   1.00 50.00 .
ATOM   773  CB  CB  . ALA ALA . . 107 107 . 8.316   25.341 7.415   1.00 50.00 .
ATOM   774  N   N   . PHE PHE . . 108 108 . 8.278   21.745 7.276   1.00 50.00 .
ATOM   775  CA  CA  . PHE PHE . . 108 108 . 8.737   20.471 7.855   1.00 50.00 .
ATOM   776  C   C   . PHE PHE . . 108 108 . 10.034  19.915 7.255   1.00 50.00 .
ATOM   777  O   O   . PHE PHE . . 108 108 . 10.365  20.158 6.097   1.00 50.00 .
ATOM   778  CB  CB  . PHE PHE . . 108 108 . 7.590   19.458 7.753   1.00 50.00 .
ATOM   779  CG  CG  . PHE PHE . . 108 108 . 8.011   18.001 7.960   1.00 50.00 .
ATOM   780  CD1 CD1 . PHE PHE . . 108 108 . 8.154   17.507 9.248   1.00 50.00 .
ATOM   781  CD2 CD2 . PHE PHE . . 108 108 . 8.388   17.232 6.865   1.00 50.00 .
ATOM   782  CE1 CE1 . PHE PHE . . 108 108 . 8.659   16.230 9.446   1.00 50.00 .
ATOM   783  CE2 CE2 . PHE PHE . . 108 108 . 8.899   15.956 7.062   1.00 50.00 .
ATOM   784  CZ  CZ  . PHE PHE . . 108 108 . 9.028   15.454 8.352   1.00 50.00 .
ATOM   785  N   N   . GLY GLY . . 109 109 . 10.700  19.136 8.117   1.00 50.00 .
ATOM   786  CA  CA  . GLY GLY . . 109 109 . 11.834  18.269 7.759   1.00 50.00 .
ATOM   787  C   C   . GLY GLY . . 109 109 . 12.974  18.991 7.039   1.00 50.00 .
ATOM   788  O   O   . GLY GLY . . 109 109 . 13.541  19.967 7.505   1.00 50.00 .
ATOM   789  N   N   . GLU GLU . . 110 110 . 13.444  18.326 6.006   1.00 50.00 .
ATOM   790  CA  CA  . GLU GLU . . 110 110 . 14.555  18.797 5.155   1.00 50.00 .
ATOM   791  C   C   . GLU GLU . . 110 110 . 14.558  17.912 3.901   1.00 50.00 .
ATOM   792  O   O   . GLU GLU . . 110 110 . 13.734  17.002 3.814   1.00 50.00 .
ATOM   793  CB  CB  . GLU GLU . . 110 110 . 15.888  18.722 5.930   1.00 50.00 .
ATOM   794  CG  CG  . GLU GLU . . 110 110 . 16.184  17.326 6.491   1.00 50.00 .
ATOM   795  CD  CD  . GLU GLU . . 110 110 . 17.479  17.284 7.297   1.00 50.00 .
ATOM   796  OE1 OE1 . GLU GLU . . 110 110 . 17.527  17.941 8.357   1.00 50.00 .
ATOM   797  OE2 OE2 . GLU GLU . . 110 110 . 18.383  16.567 6.816   1.00 50.00 .
ATOM   798  N   N   . ASP ASP . . 111 111 . 15.647  17.946 3.140   1.00 50.00 .
ATOM   799  CA  CA  . ASP ASP . . 111 111 . 15.767  17.059 1.962   1.00 50.00 .
ATOM   800  C   C   . ASP ASP . . 111 111 . 15.835  15.573 2.345   1.00 50.00 .
ATOM   801  O   O   . ASP ASP . . 111 111 . 15.251  14.732 1.670   1.00 50.00 .
ATOM   802  CB  CB  . ASP ASP . . 111 111 . 16.960  17.440 1.082   1.00 50.00 .
ATOM   803  CG  CG  . ASP ASP . . 111 111 . 16.817  18.824 0.435   1.00 50.00 .
ATOM   804  OD1 OD1 . ASP ASP . . 111 111 . 15.690  19.368 0.433   1.00 50.00 .
ATOM   805  OD2 OD2 . ASP ASP . . 111 111 . 17.878  19.334 0.018   1.00 50.00 .
ATOM   806  N   N   . GLY GLY . . 112 112 . 16.436  15.323 3.518   1.00 50.00 .
ATOM   807  CA  CA  . GLY GLY . . 112 112 . 16.576  13.969 4.088   1.00 50.00 .
ATOM   808  C   C   . GLY GLY . . 112 112 . 15.256  13.400 4.633   1.00 50.00 .
ATOM   809  O   O   . GLY GLY . . 112 112 . 14.966  12.219 4.473   1.00 50.00 .
ATOM   810  N   N   . GLU GLU . . 113 113 . 14.463  14.276 5.244   1.00 50.00 .
ATOM   811  CA  CA  . GLU GLU . . 113 113 . 13.210  13.897 5.931   1.00 50.00 .
ATOM   812  C   C   . GLU GLU . . 113 113 . 11.953  13.992 5.051   1.00 50.00 .
ATOM   813  O   O   . GLU GLU . . 113 113 . 10.889  13.508 5.438   1.00 50.00 .
ATOM   814  CB  CB  . GLU GLU . . 113 113 . 13.019  14.779 7.166   1.00 50.00 .
ATOM   815  CG  CG  . GLU GLU . . 113 113 . 14.136  14.590 8.195   1.00 50.00 .
ATOM   816  CD  CD  . GLU GLU . . 113 113 . 13.995  15.611 9.323   1.00 50.00 .
ATOM   817  OE1 OE1 . GLU GLU . . 113 113 . 13.280  15.290 10.294  1.00 50.00 .
ATOM   818  OE2 OE2 . GLU GLU . . 113 113 . 14.567  16.710 9.152   1.00 50.00 .
ATOM   819  N   N   . GLY GLY . . 114 114 . 12.077  14.719 3.931   1.00 50.00 .
ATOM   820  CA  CA  . GLY GLY . . 114 114 . 10.957  15.009 3.021   1.00 50.00 .
ATOM   821  C   C   . GLY GLY . . 114 114 . 10.211  13.739 2.599   1.00 50.00 .
ATOM   822  O   O   . GLY GLY . . 114 114 . 10.803  12.702 2.314   1.00 50.00 .
ATOM   823  N   N   . ASP ASP . . 115 115 . 8.893   13.829 2.696   1.00 50.00 .
ATOM   824  CA  CA  . ASP ASP . . 115 115 . 7.990   12.737 2.285   1.00 50.00 .
ATOM   825  C   C   . ASP ASP . . 115 115 . 6.916   13.252 1.307   1.00 50.00 .
ATOM   826  O   O   . ASP ASP . . 115 115 . 6.877   14.443 0.993   1.00 50.00 .
ATOM   827  CB  CB  . ASP ASP . . 115 115 . 7.406   12.073 3.547   1.00 50.00 .
ATOM   828  CG  CG  . ASP ASP . . 115 115 . 6.606   13.025 4.447   1.00 50.00 .
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ATOM   829  OD1 OD1 . ASP ASP . . 115 115 . 5.980   13.950 3.878   1.00 50.00 .
ATOM   830  OD2 OD2 . ASP ASP . . 115 115 . 6.589   12.791 5.667   1.00 50.00 .
ATOM   831  N   N   . ASP ASP . . 116 116 . 5.908   12.414 1.059   1.00 50.00 .
ATOM   832  CA  CA  . ASP ASP . . 116 116 . 4.784   12.754 0.158   1.00 50.00 .
ATOM   833  C   C   . ASP ASP . . 116 116 . 3.997   13.997 0.610   1.00 50.00 .
ATOM   834  O   O   . ASP ASP . . 116 116 . 3.587   14.807 -0.216  1.00 50.00 .
ATOM   835  CB  CB  . ASP ASP . . 116 116 . 3.829   11.564 0.019   1.00 50.00 .
ATOM   836  CG  CG  . ASP ASP . . 116 116 . 2.778   11.731 -1.094  1.00 50.00 .
ATOM   837  OD1 OD1 . ASP ASP . . 116 116 . 2.338   12.868 -1.347  1.00 50.00 .
ATOM   838  OD2 OD2 . ASP ASP . . 116 116 . 2.439   10.693 -1.690  1.00 50.00 .
ATOM   839  N   N   . LEU LEU . . 117 117 . 3.906   14.213 1.915   1.00 50.00 .
ATOM   840  CA  CA  . LEU LEU . . 117 117 . 3.176   15.371 2.462   1.00 50.00 .
ATOM   841  C   C   . LEU LEU . . 117 117 . 3.818   16.733 2.104   1.00 50.00 .
ATOM   842  O   O   . LEU LEU . . 117 117 . 3.348   17.791 2.513   1.00 50.00 .
ATOM   843  CB  CB  . LEU LEU . . 117 117 . 3.045   15.230 3.978   1.00 50.00 .
ATOM   844  CG  CG  . LEU LEU . . 117 117 . 2.152   14.146 4.592   1.00 50.00 .
ATOM   845  CD1 CD1 . LEU LEU . . 117 117 . 2.619   12.724 4.274   1.00 50.00 .
ATOM   846  CD2 CD2 . LEU LEU . . 117 117 . 2.150   14.319 6.112   1.00 50.00 .
ATOM   847  N   N   . ASP ASP . . 118 118 . 4.986   16.696 1.464   1.00 50.00 .
ATOM   848  CA  CA  . ASP ASP . . 118 118 . 5.628   17.908 0.923   1.00 50.00 .
ATOM   849  C   C   . ASP ASP . . 118 118 . 5.171   18.285 -0.493  1.00 50.00 .
ATOM   850  O   O   . ASP ASP . . 118 118 . 5.538   19.348 -0.986  1.00 50.00 .
ATOM   851  CB  CB  . ASP ASP . . 118 118 . 7.153   17.775 0.961   1.00 50.00 .
ATOM   852  CG  CG  . ASP ASP . . 118 118 . 7.704   17.726 2.385   1.00 50.00 .
ATOM   853  OD1 OD1 . ASP ASP . . 118 118 . 7.069   18.308 3.287   1.00 50.00 .
ATOM   854  OD2 OD2 . ASP ASP . . 118 118 . 8.710   17.017 2.580   1.00 50.00 .
ATOM   855  N   N   . LEU LEU . . 119 119 . 4.513   17.357 -1.183  1.00 50.00 .
ATOM   856  CA  CA  . LEU LEU . . 119 119 . 4.108   17.578 -2.582  1.00 50.00 .
ATOM   857  C   C   . LEU LEU . . 119 119 . 2.854   18.448 -2.686  1.00 50.00 .
ATOM   858  O   O   . LEU LEU . . 119 119 . 1.839   18.245 -2.015  1.00 50.00 .
ATOM   859  CB  CB  . LEU LEU . . 119 119 . 3.874   16.273 -3.352  1.00 50.00 .
ATOM   860  CG  CG  . LEU LEU . . 119 119 . 5.074   15.423 -3.748  1.00 50.00 .
ATOM   861  CD1 CD1 . LEU LEU . . 119 119 . 5.861   15.037 -2.507  1.00 50.00 .
ATOM   862  CD2 CD2 . LEU LEU . . 119 119 . 4.592   14.158 -4.463  1.00 50.00 .
ATOM   863  N   N   . TRP TRP . . 120 120 . 2.964   19.373 -3.622  1.00 50.00 .
ATOM   864  CA  CA  . TRP TRP . . 120 120 . 1.875   20.294 -3.981  1.00 50.00 .
ATOM   865  C   C   . TRP TRP . . 120 120 . 1.718   20.283 -5.500  1.00 50.00 .
ATOM   866  O   O   . TRP TRP . . 120 120 . 2.705   20.357 -6.230  1.00 50.00 .
ATOM   867  CB  CB  . TRP TRP . . 120 120 . 2.217   21.719 -3.532  1.00 50.00 .
ATOM   868  CG  CG  . TRP TRP . . 120 120 . 2.394   21.826 -2.018  1.00 50.00 .
ATOM   869  CD1 CD1 . TRP TRP . . 120 120 . 3.509   21.556 -1.344  1.00 50.00 .
ATOM   870  CD2 CD2 . TRP TRP . . 120 120 . 1.435   22.212 -1.098  1.00 50.00 .
ATOM   871  NE1 NE1 . TRP TRP . . 120 120 . 3.307   21.748 -0.043  1.00 50.00 .
ATOM   872  CE2 CE2 . TRP TRP . . 120 120 . 2.055   22.160 0.138   1.00 50.00 .
ATOM   873  CE3 CE3 . TRP TRP . . 120 120 . 0.126   22.662 -1.219  1.00 50.00 .
ATOM   874  CZ2 CZ2 . TRP TRP . . 120 120 . 1.365   22.575 1.262   1.00 50.00 .
ATOM   875  CZ3 CZ3 . TRP TRP . . 120 120 . -0.568  23.076 -0.090  1.00 50.00 .
ATOM   876  CH2 CH2 . TRP TRP . . 120 120 . 0.057   23.037 1.149   1.00 50.00 .
ATOM   877  N   N   . THR THR . . 121 121 . 0.473   20.202 -5.942  1.00 50.00 .
ATOM   878  CA  CA  . THR THR . . 121 121 . 0.136   20.250 -7.376  1.00 50.00 .
ATOM   879  C   C   . THR THR . . 121 121 . -0.016  21.722 -7.767  1.00 50.00 .
ATOM   880  O   O   . THR THR . . 121 121 . -0.797  22.459 -7.160  1.00 50.00 .
ATOM   881  CB  CB  . THR THR . . 121 121 . -1.173  19.494 -7.657  1.00 50.00 .
ATOM   882  OG1 OG1 . THR THR . . 121 121 . -1.050  18.146 -7.197  1.00 50.00 .
ATOM   883  CG2 CG2 . THR THR . . 121 121 . -1.545  19.497 -9.145  1.00 50.00 .
ATOM   884  N   N   . VAL VAL . . 122 122 . 0.736   22.099 -8.794  1.00 50.00 .
ATOM   885  CA  CA  . VAL VAL . . 122 122 . 0.670   23.459 -9.362  1.00 50.00 .
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ATOM   886  C   C   . VAL VAL . . 122 122 . -0.569  23.536 -10.268 1.00 50.00 .
ATOM   887  O   O   . VAL VAL . . 122 122 . -0.653  22.839 -11.279 1.00 50.00 .
ATOM   888  CB  CB  . VAL VAL . . 122 122 . 1.944   23.796 -10.161 1.00 50.00 .
ATOM   889  CG1 CG1 . VAL VAL . . 122 122 . 1.947   25.257 -10.619 1.00 50.00 .
ATOM   890  CG2 CG2 . VAL VAL . . 122 122 . 3.219   23.545 -9.354  1.00 50.00 .
ATOM   891  N   N   . ARG ARG . . 123 123 . -1.477  24.437 -9.918  1.00 50.00 .
ATOM   892  CA  CA  . ARG ARG . . 123 123 . -2.715  24.640 -10.691 1.00 50.00 .
ATOM   893  C   C   . ARG ARG . . 123 123 . -2.682  25.940 -11.496 1.00 50.00 .
ATOM   894  O   O   . ARG ARG . . 123 123 . -3.079  27.008 -11.024 1.00 50.00 .
ATOM   895  CB  CB  . ARG ARG . . 123 123 . -3.922  24.667 -9.762  1.00 50.00 .
ATOM   896  CG  CG  . ARG ARG . . 123 123 . -4.136  23.321 -9.084  1.00 50.00 .
ATOM   897  CD  CD  . ARG ARG . . 123 123 . -5.348  23.444 -8.175  1.00 50.00 .
ATOM   898  NE  NE  . ARG ARG . . 123 123 . -6.598  23.699 -8.924  1.00 50.00 .
ATOM   899  CZ  CZ  . ARG ARG . . 123 123 . -7.359  22.771 -9.511  1.00 50.00 .
ATOM   900  NH1 NH1 . ARG ARG . . 123 123 . -7.040  21.484 -9.446  1.00 50.00 .
ATOM   901  NH2 NH2 . ARG ARG . . 123 123 . -8.445  23.119 -10.187 1.00 50.00 .
ATOM   902  N   N   . CYS CYS . . 124 124 . -2.082  25.823 -12.672 1.00 50.00 .
ATOM   903  CA  CA  . CYS CYS . . 124 124 . -2.046  26.920 -13.653 1.00 50.00 .
ATOM   904  C   C   . CYS CYS . . 124 124 . -2.669  26.498 -14.988 1.00 50.00 .
ATOM   905  O   O   . CYS CYS . . 124 124 . -2.792  25.310 -15.281 1.00 50.00 .
ATOM   906  CB  CB  . CYS CYS . . 124 124 . -0.624  27.475 -13.820 1.00 50.00 .
ATOM   907  SG  SG  . CYS CYS . . 124 124 . 0.724   26.273 -14.126 1.00 50.00 .
ATOM   908  N   N   . SER SER . . 125 125 . -3.178  27.488 -15.706 1.00 50.00 .
ATOM   909  CA  CA  . SER SER . . 125 125 . -3.803  27.276 -17.027 1.00 50.00 .
ATOM   910  C   C   . SER SER . . 125 125 . -2.743  27.057 -18.110 1.00 50.00 .
ATOM   911  O   O   . SER SER . . 125 125 . -1.630  27.580 -18.029 1.00 50.00 .
ATOM   912  CB  CB  . SER SER . . 125 125 . -4.700  28.456 -17.403 1.00 50.00 .
ATOM   913  OG  OG  . SER SER . . 125 125 . -5.761  28.552 -16.450 1.00 50.00 .
ATOM   914  N   N   . GLY GLY . . 126 126 . -3.141  26.287 -19.128 1.00 50.00 .
ATOM   915  CA  CA  . GLY GLY . . 126 126 . -2.208  25.841 -20.180 1.00 50.00 .
ATOM   916  C   C   . GLY GLY . . 126 126 . -1.390  24.676 -19.613 1.00 50.00 .
ATOM   917  O   O   . GLY GLY . . 126 126 . -1.897  23.883 -18.821 1.00 50.00 .
ATOM   918  N   N   . GLN GLN . . 127 127 . -0.120  24.612 -19.989 1.00 50.00 .
ATOM   919  CA  CA  . GLN GLN . . 127 127 . 0.751   23.521 -19.508 1.00 50.00 .
ATOM   920  C   C   . GLN GLN . . 127 127 . 2.226   23.924 -19.448 1.00 50.00 .
ATOM   921  O   O   . GLN GLN . . 127 127 . 3.078   23.401 -20.164 1.00 50.00 .
ATOM   922  CB  CB  . GLN GLN . . 127 127 . 0.547   22.271 -20.370 1.00 50.00 .
ATOM   923  CG  CG  . GLN GLN . . 127 127 . 1.299   21.087 -19.761 1.00 50.00 .
ATOM   924  CD  CD  . GLN GLN . . 127 127 . 1.157   19.839 -20.613 1.00 50.00 .
ATOM   925  OE1 OE1 . GLN GLN . . 127 127 . 1.475   19.776 -21.792 1.00 50.00 .
ATOM   926  NE2 NE2 . GLN GLN . . 127 127 . 0.651   18.817 -19.969 1.00 50.00 .
ATOM   927  N   N   . HIS HIS . . 128 128 . 2.471   24.942 -18.635 1.00 50.00 .
ATOM   928  CA  CA  . HIS HIS . . 128 128 . 3.827   25.480 -18.444 1.00 50.00 .
ATOM   929  C   C   . HIS HIS . . 128 128 . 3.711   26.677 -17.510 1.00 50.00 .
ATOM   930  O   O   . HIS HIS . . 128 128 . 3.172   27.691 -17.912 1.00 50.00 .
ATOM   931  CB  CB  . HIS HIS . . 128 128 . 4.453   25.868 -19.803 1.00 50.00 .
ATOM   932  CG  CG  . HIS HIS . . 128 128 . 3.703   26.874 -20.689 1.00 50.00 .
ATOM   933  ND1 ND1 . HIS HIS . . 128 128 . 2.540   26.678 -21.287 1.00 50.00 .
ATOM   934  CD2 CD2 . HIS HIS . . 128 128 . 4.153   28.049 -21.146 1.00 50.00 .
ATOM   935  CE1 CE1 . HIS HIS . . 128 128 . 2.258   27.739 -22.044 1.00 50.00 .
ATOM   936  NE2 NE2 . HIS HIS . . 128 128 . 3.271   28.597 -21.957 1.00 50.00 .
ATOM   937  N   N   . TRP TRP . . 129 129 . 3.960   26.491 -16.230 1.00 50.00 .
ATOM   938  CA  CA  . TRP TRP . . 129 129 . 3.842   27.553 -15.219 1.00 50.00 .
ATOM   939  C   C   . TRP TRP . . 129 129 . 4.607   28.833 -15.597 1.00 50.00 .
ATOM   940  O   O   . TRP TRP . . 129 129 . 5.832   28.853 -15.612 1.00 50.00 .
ATOM   941  CB  CB  . TRP TRP . . 129 129 . 4.372   26.970 -13.913 1.00 50.00 .
ATOM   942  CG  CG  . TRP TRP . . 129 129 . 4.269   27.878 -12.692 1.00 50.00 .
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ATOM   943  CD1 CD1 . TRP TRP . . 129 129 . 3.570   29.007 -12.619 1.00 50.00 .
ATOM   944  CD2 CD2 . TRP TRP . . 129 129 . 4.726   27.582 -11.413 1.00 50.00 .
ATOM   945  NE1 NE1 . TRP TRP . . 129 129 . 3.549   29.433 -11.363 1.00 50.00 .
ATOM   946  CE2 CE2 . TRP TRP . . 129 129 . 4.244   28.591 -10.605 1.00 50.00 .
ATOM   947  CE3 CE3 . TRP TRP . . 129 129 . 5.496   26.573 -10.846 1.00 50.00 .
ATOM   948  CZ2 CZ2 . TRP TRP . . 129 129 . 4.545   28.582 -9.254  1.00 50.00 .
ATOM   949  CZ3 CZ3 . TRP TRP . . 129 129 . 5.822   26.571 -9.504  1.00 50.00 .
ATOM   950  CH2 CH2 . TRP TRP . . 129 129 . 5.335   27.583 -8.694  1.00 50.00 .
ATOM   951  N   N   . GLU GLU . . 130 130 . 3.862   29.870 -15.962 1.00 50.00 .
ATOM   952  CA  CA  . GLU GLU . . 130 130 . 4.427   31.160 -16.401 1.00 50.00 .
ATOM   953  C   C   . GLU GLU . . 130 130 . 4.502   32.072 -15.180 1.00 50.00 .
ATOM   954  O   O   . GLU GLU . . 130 130 . 3.547   32.172 -14.408 1.00 50.00 .
ATOM   955  CB  CB  . GLU GLU . . 130 130 . 3.603   31.886 -17.483 1.00 50.00 .
ATOM   956  CG  CG  . GLU GLU . . 130 130 . 3.322   31.093 -18.770 1.00 50.00 .
ATOM   957  CD  CD  . GLU GLU . . 130 130 . 2.236   30.018 -18.601 1.00 50.00 .
ATOM   958  OE1 OE1 . GLU GLU . . 130 130 . 1.649   29.900 -17.506 1.00 50.00 .
ATOM   959  OE2 OE2 . GLU GLU . . 130 130 . 1.953   29.314 -19.586 1.00 50.00 .
ATOM   960  N   N   . ARG ARG . . 131 131 . 5.592   32.826 -15.128 1.00 50.00 .
ATOM   961  CA  CA  . ARG ARG . . 131 131 . 5.839   33.786 -14.033 1.00 50.00 .
ATOM   962  C   C   . ARG ARG . . 131 131 . 4.807   34.932 -13.981 1.00 50.00 .
ATOM   963  O   O   . ARG ARG . . 131 131 . 4.634   35.583 -12.955 1.00 50.00 .
ATOM   964  CB  CB  . ARG ARG . . 131 131 . 7.258   34.338 -14.169 1.00 50.00 .
ATOM   965  CG  CG  . ARG ARG . . 131 131 . 7.637   35.253 -12.999 1.00 50.00 .
ATOM   966  CD  CD  . ARG ARG . . 131 131 . 9.038   35.834 -13.173 1.00 50.00 .
ATOM   967  NE  NE  . ARG ARG . . 131 131 . 10.032  34.756 -13.030 1.00 50.00 .
ATOM   968  CZ  CZ  . ARG ARG . . 131 131 . 10.514  34.279 -11.880 1.00 50.00 .
ATOM   969  NH1 NH1 . ARG ARG . . 131 131 . 10.118  34.763 -10.709 1.00 50.00 .
ATOM   970  NH2 NH2 . ARG ARG . . 131 131 . 11.411  33.306 -11.896 1.00 50.00 .
ATOM   971  N   N   . GLU GLU . . 132 132 . 4.081   35.118 -15.077 1.00 50.00 .
ATOM   972  CA  CA  . GLU GLU . . 132 132 . 3.101   36.210 -15.199 1.00 50.00 .
ATOM   973  C   C   . GLU GLU . . 132 132 . 1.646   35.805 -14.924 1.00 50.00 .
ATOM   974  O   O   . GLU GLU . . 132 132 . 0.709   36.587 -15.094 1.00 50.00 .
ATOM   975  CB  CB  . GLU GLU . . 132 132 . 3.212   36.734 -16.619 1.00 50.00 .
ATOM   976  CG  CG  . GLU GLU . . 132 132 . 2.472   38.043 -16.824 1.00 50.00 .
ATOM   977  CD  CD  . GLU GLU . . 132 132 . 2.736   39.196 -15.856 1.00 50.00 .
ATOM   978  OE1 OE1 . GLU GLU . . 132 132 . 3.858   39.247 -15.313 1.00 50.00 .
ATOM   979  OE2 OE2 . GLU GLU . . 132 132 . 1.826   40.044 -15.737 1.00 50.00 .
ATOM   980  N   N   . ALA ALA . . 133 133 . 1.464   34.571 -14.495 1.00 50.00 .
ATOM   981  CA  CA  . ALA ALA . . 133 133 . 0.110   34.059 -14.272 1.00 50.00 .
ATOM   982  C   C   . ALA ALA . . 133 133 . -0.124  33.789 -12.786 1.00 50.00 .
ATOM   983  O   O   . ALA ALA . . 133 133 . 0.802   33.489 -12.032 1.00 50.00 .
ATOM   984  CB  CB  . ALA ALA . . 133 133 . -0.086  32.792 -15.106 1.00 50.00 .
ATOM   985  N   N   . ALA ALA . . 134 134 . -1.368  34.030 -12.393 1.00 50.00 .
ATOM   986  CA  CA  . ALA ALA . . 134 134 . -1.874  33.600 -11.081 1.00 50.00 .
ATOM   987  C   C   . ALA ALA . . 134 134 . -2.032  32.071 -11.098 1.00 50.00 .
ATOM   988  O   O   . ALA ALA . . 134 134 . -2.155  31.458 -12.160 1.00 50.00 .
ATOM   989  CB  CB  . ALA ALA . . 134 134 . -3.217  34.281 -10.830 1.00 50.00 .
ATOM   990  N   N   . VAL VAL . . 135 135 . -2.013  31.482 -9.911  1.00 50.00 .
ATOM   991  CA  CA  . VAL VAL . . 135 135 . -1.976  30.015 -9.735  1.00 50.00 .
ATOM   992  C   C   . VAL VAL . . 135 135 . -2.360  29.634 -8.294  1.00 50.00 .
ATOM   993  O   O   . VAL VAL . . 135 135 . -2.295  30.461 -7.384  1.00 50.00 .
ATOM   994  CB  CB  . VAL VAL . . 135 135 . -0.582  29.494 -10.172 1.00 50.00 .
ATOM   995  CG1 CG1 . VAL VAL . . 135 135 . 0.578   30.230 -9.490  1.00 50.00 .
ATOM   996  CG2 CG2 . VAL VAL . . 135 135 . -0.392  27.982 -10.026 1.00 50.00 .
ATOM   997  N   N   . ARG ARG . . 136 136 . -2.831  28.399 -8.170  1.00 50.00 .
ATOM   998  CA  CA  . ARG ARG . . 136 136 . -3.198  27.766 -6.891  1.00 50.00 .
ATOM   999  C   C   . ARG ARG . . 136 136 . -2.427  26.449 -6.683  1.00 50.00 .
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ATOM   1000 O   O   . ARG ARG . . 136 136 . -1.784  25.965 -7.608  1.00 50.00 .
ATOM   1001 CB  CB  . ARG ARG . . 136 136 . -4.672  27.417 -6.851  1.00 50.00 .
ATOM   1002 CG  CG  . ARG ARG . . 136 136 . -5.547  28.658 -7.007  1.00 50.00 .
ATOM   1003 CD  CD  . ARG ARG . . 136 136 . -6.992  28.289 -6.684  1.00 50.00 .
ATOM   1004 NE  NE  . ARG ARG . . 136 136 . -7.513  27.281 -7.618  1.00 50.00 .
ATOM   1005 CZ  CZ  . ARG ARG . . 136 136 . -8.096  27.582 -8.774  1.00 50.00 .
ATOM   1006 NH1 NH1 . ARG ARG . . 136 136 . -8.243  28.848 -9.146  1.00 50.00 .
ATOM   1007 NH2 NH2 . ARG ARG . . 136 136 . -8.554  26.619 -9.554  1.00 50.00 .
ATOM   1008 N   N   . PHE PHE . . 137 137 . -2.192  26.193 -5.407  1.00 50.00 .
ATOM   1009 CA  CA  . PHE PHE . . 137 137 . -1.315  25.070 -4.980  1.00 50.00 .
ATOM   1010 C   C   . PHE PHE . . 137 137 . -2.155  24.145 -4.119  1.00 50.00 .
ATOM   1011 O   O   . PHE PHE . . 137 137 . -2.672  24.533 -3.071  1.00 50.00 .
ATOM   1012 CB  CB  . PHE PHE . . 137 137 . -0.081  25.554 -4.204  1.00 50.00 .
ATOM   1013 CG  CG  . PHE PHE . . 137 137 . 0.914   26.405 -5.013  1.00 50.00 .
ATOM   1014 CD1 CD1 . PHE PHE . . 137 137 . 0.485   27.248 -6.024  1.00 50.00 .
ATOM   1015 CD2 CD2 . PHE PHE . . 137 137 . 2.278   26.374 -4.744  1.00 50.00 .
ATOM   1016 CE1 CE1 . PHE PHE . . 137 137 . 1.327   28.010 -6.775  1.00 50.00 .
ATOM   1017 CE2 CE2 . PHE PHE . . 137 137 . 3.145   27.171 -5.490  1.00 50.00 .
ATOM   1018 CZ  CZ  . PHE PHE . . 137 137 . 2.671   27.989 -6.504  1.00 50.00 .
ATOM   1019 N   N   . GLN GLN . . 138 138 . -2.380  22.970 -4.682  1.00 50.00 .
ATOM   1020 CA  CA  . GLN GLN . . 138 138 . -3.244  21.976 -4.040  1.00 50.00 .
ATOM   1021 C   C   . GLN GLN . . 138 138 . -2.370  20.848 -3.503  1.00 50.00 .
ATOM   1022 O   O   . GLN GLN . . 138 138 . -1.686  20.156 -4.254  1.00 50.00 .
ATOM   1023 CB  CB  . GLN GLN . . 138 138 . -4.245  21.462 -5.069  1.00 50.00 .
ATOM   1024 CG  CG  . GLN GLN . . 138 138 . -5.229  20.476 -4.442  1.00 50.00 .
ATOM   1025 CD  CD  . GLN GLN . . 138 138 . -6.251  19.945 -5.446  1.00 50.00 .
ATOM   1026 OE1 OE1 . GLN GLN . . 138 138 . -6.277  20.252 -6.628  1.00 50.00 .
ATOM   1027 NE2 NE2 . GLN GLN . . 138 138 . -7.115  19.087 -4.970  1.00 50.00 .
ATOM   1028 N   N   . HIS HIS . . 139 139 . -2.395  20.714 -2.184  1.00 50.00 .
ATOM   1029 CA  CA  . HIS HIS . . 139 139 . -1.660  19.652 -1.481  1.00 50.00 .
ATOM   1030 C   C   . HIS HIS . . 139 139 . -2.073  18.307 -2.084  1.00 50.00 .
ATOM   1031 O   O   . HIS HIS . . 139 139 . -3.244  17.944 -2.043  1.00 50.00 .
ATOM   1032 CB  CB  . HIS HIS . . 139 139 . -2.022  19.680 0.005   1.00 50.00 .
ATOM   1033 CG  CG  . HIS HIS . . 139 139 . -1.113  18.772 0.835   1.00 50.00 .
ATOM   1034 ND1 ND1 . HIS HIS . . 139 139 . -1.504  17.843 1.703   1.00 50.00 .
ATOM   1035 CD2 CD2 . HIS HIS . . 139 139 . 0.212   18.836 0.879   1.00 50.00 .
ATOM   1036 CE1 CE1 . HIS HIS . . 139 139 . -0.438  17.318 2.285   1.00 50.00 .
ATOM   1037 NE2 NE2 . HIS HIS . . 139 139 . 0.611   17.935 1.768   1.00 50.00 .
ATOM   1038 N   N   . VAL VAL . . 140 140 . -1.091  17.563 -2.568  1.00 50.00 .
ATOM   1039 CA  CA  . VAL VAL . . 140 140 . -1.346  16.280 -3.260  1.00 50.00 .
ATOM   1040 C   C   . VAL VAL . . 140 140 . -2.070  15.253 -2.360  1.00 50.00 .
ATOM   1041 O   O   . VAL VAL . . 140 140 . -3.033  14.617 -2.775  1.00 50.00 .
ATOM   1042 CB  CB  . VAL VAL . . 140 140 . -0.017  15.714 -3.795  1.00 50.00 .
ATOM   1043 CG1 CG1 . VAL VAL . . 140 140 . 0.919   15.293 -2.662  1.00 50.00 .
ATOM   1044 CG2 CG2 . VAL VAL . . 140 140 . -0.227  14.522 -4.732  1.00 50.00 .
ATOM   1045 N   N   . GLY GLY . . 141 141 . -1.568  15.155 -1.119  1.00 50.00 .
ATOM   1046 CA  CA  . GLY GLY . . 141 141 . -1.957  14.100 -0.169  1.00 50.00 .
ATOM   1047 C   C   . GLY GLY . . 141 141 . -3.386  14.291 0.329   1.00 50.00 .
ATOM   1048 O   O   . GLY GLY . . 141 141 . -4.196  13.371 0.335   1.00 50.00 .
ATOM   1049 N   N   . THR THR . . 142 142 . -3.670  15.538 0.676   1.00 50.00 .
ATOM   1050 CA  CA  . THR THR . . 142 142 . -4.960  15.888 1.296   1.00 50.00 .
ATOM   1051 C   C   . THR THR . . 142 142 . -5.963  16.523 0.333   1.00 50.00 .
ATOM   1052 O   O   . THR THR . . 142 142 . -7.144  16.530 0.659   1.00 50.00 .
ATOM   1053 CB  CB  . THR THR . . 142 142 . -4.793  16.790 2.522   1.00 50.00 .
ATOM   1054 OG1 OG1 . THR THR . . 142 142 . -4.152  18.009 2.130   1.00 50.00 .
ATOM   1055 CG2 CG2 . THR THR . . 142 142 . -4.009  16.080 3.633   1.00 50.00 .
ATOM   1056 N   N   . SER SER . . 143 143 . -5.462  17.187 -0.719  1.00 50.00 .
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ATOM   1057 CA  CA  . SER SER . . 143 143 . -6.257  17.894 -1.757  1.00 50.00 .
ATOM   1058 C   C   . SER SER . . 143 143 . -6.597  19.340 -1.362  1.00 50.00 .
ATOM   1059 O   O   . SER SER . . 143 143 . -7.178  20.123 -2.108  1.00 50.00 .
ATOM   1060 CB  CB  . SER SER . . 143 143 . -7.512  17.086 -2.127  1.00 50.00 .
ATOM   1061 OG  OG  . SER SER . . 143 143 . -8.428  17.790 -2.955  1.00 50.00 .
ATOM   1062 N   N   . VAL VAL . . 144 144 . -6.131  19.716 -0.184  1.00 50.00 .
ATOM   1063 CA  CA  . VAL VAL . . 144 144 . -6.342  21.053 0.357   1.00 50.00 .
ATOM   1064 C   C   . VAL VAL . . 144 144 . -5.580  22.098 -0.468  1.00 50.00 .
ATOM   1065 O   O   . VAL VAL . . 144 144 . -4.429  21.887 -0.844  1.00 50.00 .
ATOM   1066 CB  CB  . VAL VAL . . 144 144 . -5.898  20.942 1.810   1.00 50.00 .
ATOM   1067 CG1 CG1 . VAL VAL . . 144 144 . -4.425  21.261 2.095   1.00 50.00 .
ATOM   1068 CG2 CG2 . VAL VAL . . 144 144 . -6.875  21.752 2.624   1.00 50.00 .
ATOM   1069 N   N   . PHE PHE . . 145 145 . -6.160  23.278 -0.538  1.00 50.00 .
ATOM   1070 CA  CA  . PHE PHE . . 145 145 . -5.501  24.387 -1.231  1.00 50.00 .
ATOM   1071 C   C   . PHE PHE . . 145 145 . -4.773  25.288 -0.235  1.00 50.00 .
ATOM   1072 O   O   . PHE PHE . . 145 145 . -5.290  25.566 0.851   1.00 50.00 .
ATOM   1073 CB  CB  . PHE PHE . . 145 145 . -6.566  25.199 -1.937  1.00 50.00 .
ATOM   1074 CG  CG  . PHE PHE . . 145 145 . -7.255  24.436 -3.069  1.00 50.00 .
ATOM   1075 CD1 CD1 . PHE PHE . . 145 145 . -6.683  24.386 -4.333  1.00 50.00 .
ATOM   1076 CD2 CD2 . PHE PHE . . 145 145 . -8.478  23.821 -2.830  1.00 50.00 .
ATOM   1077 CE1 CE1 . PHE PHE . . 145 145 . -7.348  23.695 -5.334  1.00 50.00 .
ATOM   1078 CE2 CE2 . PHE PHE . . 145 145 . -9.128  23.130 -3.843  1.00 50.00 .
ATOM   1079 CZ  CZ  . PHE PHE . . 145 145 . -8.556  23.060 -5.103  1.00 50.00 .
ATOM   1080 N   N   . LEU LEU . . 146 146 . -3.567  25.676 -0.636  1.00 50.00 .
ATOM   1081 CA  CA  . LEU LEU . . 146 146 . -2.770  26.677 0.089   1.00 50.00 .
ATOM   1082 C   C   . LEU LEU . . 146 146 . -3.512  28.018 0.024   1.00 50.00 .
ATOM   1083 O   O   . LEU LEU . . 146 146 . -3.714  28.572 -1.055  1.00 50.00 .
ATOM   1084 CB  CB  . LEU LEU . . 146 146 . -1.390  26.812 -0.565  1.00 50.00 .
ATOM   1085 CG  CG  . LEU LEU . . 146 146 . -0.462  27.740 0.225   1.00 50.00 .
ATOM   1086 CD1 CD1 . LEU LEU . . 146 146 . -0.088  27.103 1.563   1.00 50.00 .
ATOM   1087 CD2 CD2 . LEU LEU . . 146 146 . 0.795   28.056 -0.585  1.00 50.00 .
ATOM   1088 N   N   . SER SER . . 147 147 . -3.908  28.499 1.194   1.00 50.00 .
ATOM   1089 CA  CA  . SER SER . . 147 147 . -4.769  29.693 1.283   1.00 50.00 .
ATOM   1090 C   C   . SER SER . . 147 147 . -4.487  30.456 2.573   1.00 50.00 .
ATOM   1091 O   O   . SER SER . . 147 147 . -4.046  29.875 3.565   1.00 50.00 .
ATOM   1092 CB  CB  . SER SER . . 147 147 . -6.249  29.287 1.282   1.00 50.00 .
ATOM   1093 OG  OG  . SER SER . . 147 147 . -6.575  28.541 0.104   1.00 50.00 .
ATOM   1094 N   N   . VAL VAL . . 148 148 . -4.606  31.773 2.493   1.00 50.00 .
ATOM   1095 CA  CA  . VAL VAL . . 148 148 . -4.622  32.597 3.718   1.00 50.00 .
ATOM   1096 C   C   . VAL VAL . . 148 148 . -6.026  32.485 4.330   1.00 50.00 .
ATOM   1097 O   O   . VAL VAL . . 148 148 . -7.004  32.246 3.636   1.00 50.00 .
ATOM   1098 CB  CB  . VAL VAL . . 148 148 . -4.090  34.030 3.474   1.00 50.00 .
ATOM   1099 CG1 CG1 . VAL VAL . . 148 148 . -4.771  34.776 2.340   1.00 50.00 .
ATOM   1100 CG2 CG2 . VAL VAL . . 148 148 . -4.166  34.909 4.722   1.00 50.00 .
ATOM   1101 N   N   . THR THR . . 149 149 . -6.062  32.305 5.637   1.00 50.00 .
ATOM   1102 CA  CA  . THR THR . . 149 149 . -7.358  32.143 6.340   1.00 50.00 .
ATOM   1103 C   C   . THR THR . . 149 149 . -8.203  33.422 6.280   1.00 50.00 .
ATOM   1104 O   O   . THR THR . . 149 149 . -9.421  33.392 6.434   1.00 50.00 .
ATOM   1105 CB  CB  . THR THR . . 149 149 . -7.159  31.793 7.811   1.00 50.00 .
ATOM   1106 OG1 OG1 . THR THR . . 149 149 . -6.453  32.856 8.437   1.00 50.00 .
ATOM   1107 CG2 CG2 . THR THR . . 149 149 . -6.331  30.536 7.995   1.00 50.00 .
ATOM   1108 N   N   . GLY GLY . . 150 150 . -7.446  34.523 6.331   1.00 50.00 .
ATOM   1109 CA  CA  . GLY GLY . . 150 150 . -7.994  35.873 6.535   1.00 50.00 .
ATOM   1110 C   C   . GLY GLY . . 150 150 . -7.703  36.399 7.938   1.00 50.00 .
ATOM   1111 O   O   . GLY GLY . . 150 150 . -7.921  37.570 8.229   1.00 50.00 .
ATOM   1112 N   N   . GLU GLU . . 151 151 . -7.201  35.505 8.777   1.00 50.00 .
ATOM   1113 CA  CA  . GLU GLU . . 151 151 . -6.802  35.826 10.144  1.00 50.00 .
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ATOM   1114 C   C   . GLU GLU . . 151 151 . -5.319  36.199 10.185  1.00 50.00 .
ATOM   1115 O   O   . GLU GLU . . 151 151 . -4.512  35.789 9.346   1.00 50.00 .
ATOM   1116 CB  CB  . GLU GLU . . 151 151 . -7.106  34.600 10.992  1.00 50.00 .
ATOM   1117 CG  CG  . GLU GLU . . 151 151 . -6.789  34.817 12.471  1.00 50.00 .
ATOM   1118 CD  CD  . GLU GLU . . 151 151 . -7.335  33.688 13.330  1.00 50.00 .
ATOM   1119 OE1 OE1 . GLU GLU . . 151 151 . -7.602  32.602 12.766  1.00 50.00 .
ATOM   1120 OE2 OE2 . GLU GLU . . 151 151 . -7.418  33.930 14.552  1.00 50.00 .
ATOM   1121 N   N   . GLN GLN . . 152 152 . -5.030  37.030 11.172  1.00 50.00 .
ATOM   1122 CA  CA  . GLN GLN . . 152 152 . -3.670  37.506 11.455  1.00 50.00 .
ATOM   1123 C   C   . GLN GLN . . 152 152 . -3.259  37.078 12.864  1.00 50.00 .
ATOM   1124 O   O   . GLN GLN . . 152 152 . -4.095  36.903 13.750  1.00 50.00 .
ATOM   1125 CB  CB  . GLN GLN . . 152 152 . -3.621  39.030 11.376  1.00 50.00 .
ATOM   1126 CG  CG  . GLN GLN . . 152 152 . -3.960  39.540 9.976   1.00 50.00 .
ATOM   1127 CD  CD  . GLN GLN . . 152 152 . -3.802  41.056 9.868   1.00 50.00 .
ATOM   1128 OE1 OE1 . GLN GLN . . 152 152 . -2.963  41.686 10.494  1.00 50.00 .
ATOM   1129 NE2 NE2 . GLN GLN . . 152 152 . -4.627  41.664 9.045   1.00 50.00 .
ATOM   1130 N   N   . TYR TYR . . 153 153 . -1.959  36.891 13.014  1.00 50.00 .
ATOM   1131 CA  CA  . TYR TYR . . 153 153 . -1.351  36.536 14.304  1.00 50.00 .
ATOM   1132 C   C   . TYR TYR . . 153 153 . -1.012  37.793 15.110  1.00 50.00 .
ATOM   1133 O   O   . TYR TYR . . 153 153 . -0.914  38.900 14.572  1.00 50.00 .
ATOM   1134 CB  CB  . TYR TYR . . 153 153 . -0.081  35.715 14.075  1.00 50.00 .
ATOM   1135 CG  CG  . TYR TYR . . 153 153 . -0.377  34.364 13.424  1.00 50.00 .
ATOM   1136 CD1 CD1 . TYR TYR . . 153 153 . -0.723  33.268 14.204  1.00 50.00 .
ATOM   1137 CD2 CD2 . TYR TYR . . 153 153 . -0.243  34.221 12.051  1.00 50.00 .
ATOM   1138 CE1 CE1 . TYR TYR . . 153 153 . -0.935  32.034 13.607  1.00 50.00 .
ATOM   1139 CE2 CE2 . TYR TYR . . 153 153 . -0.454  32.993 11.455  1.00 50.00 .
ATOM   1140 CZ  CZ  . TYR TYR . . 153 153 . -0.803  31.902 12.233  1.00 50.00 .
ATOM   1141 OH  OH  . TYR TYR . . 153 153 . -1.029  30.707 11.642  1.00 50.00 .
ATOM   1142 N   N   . GLY GLY . . 154 154 . -0.913  37.575 16.424  1.00 50.00 .
ATOM   1143 CA  CA  . GLY GLY . . 154 154 . -0.404  38.609 17.343  1.00 50.00 .
ATOM   1144 C   C   . GLY GLY . . 154 154 . 1.099   38.374 17.538  1.00 50.00 .
ATOM   1145 O   O   . GLY GLY . . 154 154 . 1.738   37.627 16.792  1.00 50.00 .
ATOM   1146 N   N   . SER SER . . 155 155 . 1.632   38.930 18.619  1.00 50.00 .
ATOM   1147 CA  CA  . SER SER . . 155 155 . 3.034   38.670 19.007  1.00 50.00 .
ATOM   1148 C   C   . SER SER . . 155 155 . 3.263   37.155 19.221  1.00 50.00 .
ATOM   1149 O   O   . SER SER . . 155 155 . 2.302   36.424 19.472  1.00 50.00 .
ATOM   1150 CB  CB  . SER SER . . 155 155 . 3.395   39.449 20.276  1.00 50.00 .
ATOM   1151 OG  OG  . SER SER . . 155 155 . 2.562   39.023 21.353  1.00 50.00 .
ATOM   1152 N   N   . PRO PRO . . 156 156 . 4.496   36.681 18.989  1.00 50.00 .
ATOM   1153 CA  CA  . PRO PRO . . 156 156 . 5.679   37.480 18.598  1.00 50.00 .
ATOM   1154 C   C   . PRO PRO . . 156 156 . 5.804   37.729 17.083  1.00 50.00 .
ATOM   1155 O   O   . PRO PRO . . 156 156 . 6.854   38.151 16.605  1.00 50.00 .
ATOM   1156 CB  CB  . PRO PRO . . 156 156 . 6.838   36.640 19.131  1.00 50.00 .
ATOM   1157 CG  CG  . PRO PRO . . 156 156 . 6.362   35.211 18.866  1.00 50.00 .
ATOM   1158 CD  CD  . PRO PRO . . 156 156 . 4.869   35.269 19.197  1.00 50.00 .
ATOM   1159 N   N   . ILE ILE . . 157 157 . 4.717   37.488 16.353  1.00 50.00 .
ATOM   1160 CA  CA  . ILE ILE . . 157 157 . 4.642   37.656 14.886  1.00 50.00 .
ATOM   1161 C   C   . ILE ILE . . 157 157 . 3.441   38.514 14.475  1.00 50.00 .
ATOM   1162 O   O   . ILE ILE . . 157 157 . 2.679   38.188 13.563  1.00 50.00 .
ATOM   1163 CB  CB  . ILE ILE . . 157 157 . 4.592   36.298 14.173  1.00 50.00 .
ATOM   1164 CG1 CG1 . ILE ILE . . 157 157 . 3.513   35.351 14.730  1.00 50.00 .
ATOM   1165 CG2 CG2 . ILE ILE . . 157 157 . 5.995   35.717 14.207  1.00 50.00 .
ATOM   1166 CD1 CD1 . ILE ILE . . 157 157 . 3.357   34.039 13.952  1.00 50.00 .
ATOM   1167 N   N   . ARG ARG . . 158 158 . 3.294   39.619 15.194  1.00 50.00 .
ATOM   1168 CA  CA  . ARG ARG . . 158 158 . 2.132   40.507 15.031  1.00 50.00 .
ATOM   1169 C   C   . ARG ARG . . 158 158 . 1.984   40.973 13.577  1.00 50.00 .
ATOM   1170 O   O   . ARG ARG . . 158 158 . 2.956   41.388 12.947  1.00 50.00 .
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ATOM   1171 CB  CB  . ARG ARG . . 158 158 . 2.309   41.712 15.951  1.00 50.00 .
ATOM   1172 CG  CG  . ARG ARG . . 158 158 . 1.142   42.696 15.823  1.00 50.00 .
ATOM   1173 CD  CD  . ARG ARG . . 158 158 . 1.387   43.937 16.676  1.00 50.00 .
ATOM   1174 NE  NE  . ARG ARG . . 158 158 . 1.368   43.552 18.098  1.00 50.00 .
ATOM   1175 CZ  CZ  . ARG ARG . . 158 158 . 0.278   43.373 18.849  1.00 50.00 .
ATOM   1176 NH1 NH1 . ARG ARG . . 158 158 . -0.940  43.547 18.349  1.00 50.00 .
ATOM   1177 NH2 NH2 . ARG ARG . . 158 158 . 0.403   43.015 20.117  1.00 50.00 .
ATOM   1178 N   N   . GLY GLY . . 159 159 . 0.739   40.864 13.105  1.00 50.00 .
ATOM   1179 CA  CA  . GLY GLY . . 159 159 . 0.368   41.350 11.763  1.00 50.00 .
ATOM   1180 C   C   . GLY GLY . . 159 159 . 0.508   40.298 10.653  1.00 50.00 .
ATOM   1181 O   O   . GLY GLY . . 159 159 . -0.024  40.480 9.560   1.00 50.00 .
ATOM   1182 N   N   . GLN GLN . . 160 160 . 1.249   39.225 10.923  1.00 50.00 .
ATOM   1183 CA  CA  . GLN GLN . . 160 160 . 1.472   38.170 9.917   1.00 50.00 .
ATOM   1184 C   C   . GLN GLN . . 160 160 . 0.187   37.366 9.706   1.00 50.00 .
ATOM   1185 O   O   . GLN GLN . . 160 160 . -0.543  37.075 10.652  1.00 50.00 .
ATOM   1186 CB  CB  . GLN GLN . . 160 160 . 2.621   37.257 10.340  1.00 50.00 .
ATOM   1187 CG  CG  . GLN GLN . . 160 160 . 3.930   38.048 10.417  1.00 50.00 .
ATOM   1188 CD  CD  . GLN GLN . . 160 160 . 5.132   37.146 10.692  1.00 50.00 .
ATOM   1189 OE1 OE1 . GLN GLN . . 160 160 . 5.201   35.998 10.274  1.00 50.00 .
ATOM   1190 NE2 NE2 . GLN GLN . . 160 160 . 6.112   37.689 11.378  1.00 50.00 .
ATOM   1191 N   N   . HIS HIS . . 161 161 . -0.096  37.070 8.447   1.00 50.00 .
ATOM   1192 CA  CA  . HIS HIS . . 161 161 . -1.323  36.350 8.070   1.00 50.00 .
ATOM   1193 C   C   . HIS HIS . . 161 161 . -1.176  34.841 8.248   1.00 50.00 .
ATOM   1194 O   O   . HIS HIS . . 161 161 . -0.136  34.253 7.950   1.00 50.00 .
ATOM   1195 CB  CB  . HIS HIS . . 161 161 . -1.699  36.663 6.624   1.00 50.00 .
ATOM   1196 CG  CG  . HIS HIS . . 161 161 . -2.151  38.114 6.475   1.00 50.00 .
ATOM   1197 ND1 ND1 . HIS HIS . . 161 161 . -3.295  38.625 6.915   1.00 50.00 .
ATOM   1198 CD2 CD2 . HIS HIS . . 161 161 . -1.521  39.060 5.791   1.00 50.00 .
ATOM   1199 CE1 CE1 . HIS HIS . . 161 161 . -3.368  39.883 6.489   1.00 50.00 .
ATOM   1200 NE2 NE2 . HIS HIS . . 161 161 . -2.274  40.153 5.787   1.00 50.00 .
ATOM   1201 N   N   . GLU GLU . . 162 162 . -2.251  34.242 8.739   1.00 50.00 .
ATOM   1202 CA  CA  . GLU GLU . . 162 162 . -2.336  32.782 8.852   1.00 50.00 .
ATOM   1203 C   C   . GLU GLU . . 162 162 . -2.446  32.139 7.469   1.00 50.00 .
ATOM   1204 O   O   . GLU GLU . . 162 162 . -3.331  32.488 6.687   1.00 50.00 .
ATOM   1205 CB  CB  . GLU GLU . . 162 162 . -3.525  32.429 9.733   1.00 50.00 .
ATOM   1206 CG  CG  . GLU GLU . . 162 162 . -3.620  30.922 9.971   1.00 50.00 .
ATOM   1207 CD  CD  . GLU GLU . . 162 162 . -4.862  30.583 10.786  1.00 50.00 .
ATOM   1208 OE1 OE1 . GLU GLU . . 162 162 . -5.828  31.368 10.778  1.00 50.00 .
ATOM   1209 OE2 OE2 . GLU GLU . . 162 162 . -4.817  29.534 11.450  1.00 50.00 .
ATOM   1210 N   N   . VAL VAL . . 163 163 . -1.641  31.099 7.303   1.00 50.00 .
ATOM   1211 CA  CA  . VAL VAL . . 163 163 . -1.651  30.273 6.086   1.00 50.00 .
ATOM   1212 C   C   . VAL VAL . . 163 163 . -2.159  28.887 6.495   1.00 50.00 .
ATOM   1213 O   O   . VAL VAL . . 163 163 . -1.678  28.281 7.451   1.00 50.00 .
ATOM   1214 CB  CB  . VAL VAL . . 163 163 . -0.234  30.187 5.496   1.00 50.00 .
ATOM   1215 CG1 CG1 . VAL VAL . . 163 163 . -0.201  29.299 4.254   1.00 50.00 .
ATOM   1216 CG2 CG2 . VAL VAL . . 163 163 . 0.279   31.567 5.081   1.00 50.00 .
ATOM   1217 N   N   . HIS HIS . . 164 164 . -3.175  28.447 5.776   1.00 50.00 .
ATOM   1218 CA  CA  . HIS HIS . . 164 164 . -3.798  27.147 6.045   1.00 50.00 .
ATOM   1219 C   C   . HIS HIS . . 164 164 . -4.044  26.378 4.739   1.00 50.00 .
ATOM   1220 O   O   . HIS HIS . . 164 164 . -3.843  26.879 3.632   1.00 50.00 .
ATOM   1221 CB  CB  . HIS HIS . . 164 164 . -5.122  27.401 6.768   1.00 50.00 .
ATOM   1222 CG  CG  . HIS HIS . . 164 164 . -6.158  27.948 5.776   1.00 50.00 .
ATOM   1223 ND1 ND1 . HIS HIS . . 164 164 . -7.204  27.312 5.296   1.00 50.00 .
ATOM   1224 CD2 CD2 . HIS HIS . . 164 164 . -6.212  29.170 5.283   1.00 50.00 .
ATOM   1225 CE1 CE1 . HIS HIS . . 164 164 . -7.959  28.200 4.655   1.00 50.00 .
ATOM   1226 NE2 NE2 . HIS HIS . . 164 164 . -7.349  29.365 4.648   1.00 50.00 .
ATOM   1227 N   N   . GLY GLY . . 165 165 . -4.720  25.256 4.958   1.00 50.00 .
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ATOM   1228 CA  CA  . GLY GLY . . 165 165 . -5.265  24.450 3.875   1.00 50.00 .
ATOM   1229 C   C   . GLY GLY . . 165 165 . -6.785  24.490 3.992   1.00 50.00 .
ATOM   1230 O   O   . GLY GLY . . 165 165 . -7.368  23.985 4.958   1.00 50.00 .
ATOM   1231 N   N   . MET MET . . 166 166 . -7.401  24.980 2.923   1.00 50.00 .
ATOM   1232 CA  CA  . MET MET . . 166 166 . -8.867  24.990 2.784   1.00 50.00 .
ATOM   1233 C   C   . MET MET . . 166 166 . -9.324  23.855 1.845   1.00 50.00 .
ATOM   1234 O   O   . MET MET . . 166 166 . -8.708  23.663 0.793   1.00 50.00 .
ATOM   1235 CB  CB  . MET MET . . 166 166 . -9.302  26.349 2.224   1.00 50.00 .
ATOM   1236 CG  CG  . MET MET . . 166 166 . -10.822 26.554 2.175   1.00 50.00 .
ATOM   1237 SD  SD  . MET MET . . 166 166 . -11.741 26.565 3.754   1.00 50.00 .
ATOM   1238 CE  CE  . MET MET . . 166 166 . -11.369 28.219 4.290   1.00 50.00 .
ATOM   1239 N   N   . PRO PRO . . 167 167 . -10.403 23.141 2.211   1.00 50.00 .
ATOM   1240 CA  CA  . PRO PRO . . 167 167 . -11.009 22.095 1.365   1.00 50.00 .
ATOM   1241 C   C   . PRO PRO . . 167 167 . -11.469 22.636 0.002   1.00 50.00 .
ATOM   1242 O   O   . PRO PRO . . 167 167 . -11.404 21.937 -1.003  1.00 50.00 .
ATOM   1243 CB  CB  . PRO PRO . . 167 167 . -12.187 21.560 2.188   1.00 50.00 .
ATOM   1244 CG  CG  . PRO PRO . . 167 167 . -12.561 22.734 3.091   1.00 50.00 .
ATOM   1245 CD  CD  . PRO PRO . . 167 167 . -11.185 23.286 3.457   1.00 50.00 .
ATOM   1246 N   N   . SER SER . . 168 168 . -11.931 23.884 -0.009  1.00 50.00 .
ATOM   1247 CA  CA  . SER SER . . 168 168 . -12.390 24.563 -1.236  1.00 50.00 .
ATOM   1248 C   C   . SER SER . . 168 168 . -11.559 25.815 -1.538  1.00 50.00 .
ATOM   1249 O   O   . SER SER . . 168 168 . -10.851 26.343 -0.686  1.00 50.00 .
ATOM   1250 CB  CB  . SER SER . . 168 168 . -13.867 24.938 -1.093  1.00 50.00 .
ATOM   1251 OG  OG  . SER SER . . 168 168 . -14.629 23.743 -0.904  1.00 50.00 .
ATOM   1252 N   N   . ALA ALA . . 169 169 . -11.548 26.206 -2.805  1.00 50.00 .
ATOM   1253 CA  CA  . ALA ALA . . 169 169 . -10.868 27.441 -3.239  1.00 50.00 .
ATOM   1254 C   C   . ALA ALA . . 169 169 . -11.591 28.671 -2.668  1.00 50.00 .
ATOM   1255 O   O   . ALA ALA . . 169 169 . -12.813 28.778 -2.747  1.00 50.00 .
ATOM   1256 CB  CB  . ALA ALA . . 169 169 . -10.851 27.498 -4.767  1.00 50.00 .
ATOM   1257 N   N   . ASN ASN . . 170 170 . -10.816 29.532 -2.020  1.00 50.00 .
ATOM   1258 CA  CA  . ASN ASN . . 170 170 . -11.350 30.773 -1.429  1.00 50.00 .
ATOM   1259 C   C   . ASN ASN . . 170 170 . -10.729 32.003 -2.118  1.00 50.00 .
ATOM   1260 O   O   . ASN ASN . . 170 170 . -9.816  31.878 -2.937  1.00 50.00 .
ATOM   1261 CB  CB  . ASN ASN . . 170 170 . -11.094 30.728 0.089   1.00 50.00 .
ATOM   1262 CG  CG  . ASN ASN . . 170 170 . -11.735 31.908 0.830   1.00 50.00 .
ATOM   1263 OD1 OD1 . ASN ASN . . 170 170 . -12.892 32.251 0.638   1.00 50.00 .
ATOM   1264 ND2 ND2 . ASN ASN . . 170 170 . -10.902 32.721 1.438   1.00 50.00 .
ATOM   1265 N   N   . THR THR . . 171 171 . -11.086 33.179 -1.613  1.00 50.00 .
ATOM   1266 CA  CA  . THR THR . . 171 171 . -10.538 34.481 -2.059  1.00 50.00 .
ATOM   1267 C   C   . THR THR . . 171 171 . -9.063  34.664 -1.659  1.00 50.00 .
ATOM   1268 O   O   . THR THR . . 171 171 . -8.401  35.642 -1.998  1.00 50.00 .
ATOM   1269 CB  CB  . THR THR . . 171 171 . -11.367 35.635 -1.477  1.00 50.00 .
ATOM   1270 OG1 OG1 . THR THR . . 171 171 . -11.331 35.598 -0.046  1.00 50.00 .
ATOM   1271 CG2 CG2 . THR THR . . 171 171 . -12.816 35.580 -1.971  1.00 50.00 .
ATOM   1272 N   N   . HIS HIS . . 172 172 . -8.560  33.664 -0.948  1.00 50.00 .
ATOM   1273 CA  CA  . HIS HIS . . 172 172 . -7.216  33.640 -0.374  1.00 50.00 .
ATOM   1274 C   C   . HIS HIS . . 172 172 . -6.327  32.535 -0.954  1.00 50.00 .
ATOM   1275 O   O   . HIS HIS . . 172 172 . -5.185  32.374 -0.525  1.00 50.00 .
ATOM   1276 CB  CB  . HIS HIS . . 172 172 . -7.504  33.384 1.095   1.00 50.00 .
ATOM   1277 CG  CG  . HIS HIS . . 172 172 . -8.083  34.584 1.860   1.00 50.00 .
ATOM   1278 ND1 ND1 . HIS HIS . . 172 172 . -8.300  34.640 3.170   1.00 50.00 .
ATOM   1279 CD2 CD2 . HIS HIS . . 172 172 . -8.530  35.729 1.354   1.00 50.00 .
ATOM   1280 CE1 CE1 . HIS HIS . . 172 172 . -8.880  35.796 3.448   1.00 50.00 .
ATOM   1281 NE2 NE2 . HIS HIS . . 172 172 . -9.020  36.478 2.327   1.00 50.00 .
ATOM   1282 N   N   . ASN ASN . . 173 173 . -6.844  31.873 -1.986  1.00 50.00 .
ATOM   1283 CA  CA  . ASN ASN . . 173 173 . -6.150  30.747 -2.631  1.00 50.00 .
ATOM   1284 C   C   . ASN ASN . . 173 173 . -5.264  31.175 -3.811  1.00 50.00 .
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ATOM   1285 O   O   . ASN ASN . . 173 173 . -4.317  30.487 -4.184  1.00 50.00 .
ATOM   1286 CB  CB  . ASN ASN . . 173 173 . -7.187  29.719 -3.082  1.00 50.00 .
ATOM   1287 CG  CG  . ASN ASN . . 173 173 . -6.572  28.415 -3.559  1.00 50.00 .
ATOM   1288 OD1 OD1 . ASN ASN . . 173 173 . -5.391  28.246 -3.919  1.00 50.00 .
ATOM   1289 ND2 ND2 . ASN ASN . . 173 173 . -7.345  27.392 -3.657  1.00 50.00 .
ATOM   1290 N   N   . THR THR . . 174 174 . -5.518  32.341 -4.384  1.00 50.00 .
ATOM   1291 CA  CA  . THR THR . . 174 174 . -4.762  32.753 -5.578  1.00 50.00 .
ATOM   1292 C   C   . THR THR . . 174 174 . -3.397  33.328 -5.168  1.00 50.00 .
ATOM   1293 O   O   . THR THR . . 174 174 . -3.308  34.249 -4.350  1.00 50.00 .
ATOM   1294 CB  CB  . THR THR . . 174 174 . -5.585  33.739 -6.415  1.00 50.00 .
ATOM   1295 OG1 OG1 . THR THR . . 174 174 . -6.893  33.221 -6.651  1.00 50.00 .
ATOM   1296 CG2 CG2 . THR THR . . 174 174 . -4.951  33.967 -7.782  1.00 50.00 .
ATOM   1297 N   N   . TRP TRP . . 175 175 . -2.363  32.791 -5.806  1.00 50.00 .
ATOM   1298 CA  CA  . TRP TRP . . 175 175 . -0.967  33.220 -5.604  1.00 50.00 .
ATOM   1299 C   C   . TRP TRP . . 175 175 . -0.329  33.643 -6.931  1.00 50.00 .
ATOM   1300 O   O   . TRP TRP . . 175 175 . -0.800  33.298 -8.015  1.00 50.00 .
ATOM   1301 CB  CB  . TRP TRP . . 175 175 . -0.130  32.093 -4.992  1.00 50.00 .
ATOM   1302 CG  CG  . TRP TRP . . 175 175 . -0.675  31.639 -3.639  1.00 50.00 .
ATOM   1303 CD1 CD1 . TRP TRP . . 175 175 . -1.555  30.660 -3.462  1.00 50.00 .
ATOM   1304 CD2 CD2 . TRP TRP . . 175 175 . -0.331  32.119 -2.386  1.00 50.00 .
ATOM   1305 NE1 NE1 . TRP TRP . . 175 175 . -1.788  30.493 -2.162  1.00 50.00 .
ATOM   1306 CE2 CE2 . TRP TRP . . 175 175 . -1.049  31.362 -1.473  1.00 50.00 .
ATOM   1307 CE3 CE3 . TRP TRP . . 175 175 . 0.543   33.107 -1.952  1.00 50.00 .
ATOM   1308 CZ2 CZ2 . TRP TRP . . 175 175 . -0.887  31.587 -0.112  1.00 50.00 .
ATOM   1309 CZ3 CZ3 . TRP TRP . . 175 175 . 0.703   33.335 -0.591  1.00 50.00 .
ATOM   1310 CH2 CH2 . TRP TRP . . 175 175 . -0.013  32.576 0.328   1.00 50.00 .
ATOM   1311 N   N   . LYS LYS . . 176 176 . 0.684   34.482 -6.790  1.00 50.00 .
ATOM   1312 CA  CA  . LYS LYS . . 176 176 . 1.491   35.016 -7.894  1.00 50.00 .
ATOM   1313 C   C   . LYS LYS . . 176 176 . 2.972   34.877 -7.520  1.00 50.00 .
ATOM   1314 O   O   . LYS LYS . . 176 176 . 3.366   35.117 -6.379  1.00 50.00 .
ATOM   1315 CB  CB  . LYS LYS . . 176 176 . 1.099   36.483 -8.069  1.00 50.00 .
ATOM   1316 CG  CG  . LYS LYS . . 176 176 . 1.930   37.239 -9.115  1.00 50.00 .
ATOM   1317 CD  CD  . LYS LYS . . 176 176 . 1.653   38.744 -9.072  1.00 50.00 .
ATOM   1318 CE  CE  . LYS LYS . . 176 176 . 1.960   39.332 -7.691  1.00 50.00 .
ATOM   1319 NZ  NZ  . LYS LYS . . 176 176 . 1.674   40.767 -7.638  1.00 50.00 .
ATOM   1320 N   N   . ALA ALA . . 177 177 . 3.768   34.516 -8.516  1.00 50.00 .
ATOM   1321 CA  CA  . ALA ALA . . 177 177 . 5.224   34.379 -8.345  1.00 50.00 .
ATOM   1322 C   C   . ALA ALA . . 177 177 . 5.945   35.680 -8.720  1.00 50.00 .
ATOM   1323 O   O   . ALA ALA . . 177 177 . 5.480   36.466 -9.544  1.00 50.00 .
ATOM   1324 CB  CB  . ALA ALA . . 177 177 . 5.733   33.223 -9.208  1.00 50.00 .
ATOM   1325 N   N   . MET MET . . 178 178 . 7.057   35.902 -8.033  1.00 50.00 .
ATOM   1326 CA  CA  . MET MET . . 178 178 . 7.941   37.056 -8.273  1.00 50.00 .
ATOM   1327 C   C   . MET MET . . 178 178 . 9.388   36.638 -7.986  1.00 50.00 .
ATOM   1328 O   O   . MET MET . . 178 178 . 9.656   35.872 -7.056  1.00 50.00 .
ATOM   1329 CB  CB  . MET MET . . 178 178 . 7.537   38.247 -7.388  1.00 50.00 .
ATOM   1330 CG  CG  . MET MET . . 178 178 . 7.749   37.945 -5.901  1.00 50.00 .
ATOM   1331 SD  SD  . MET MET . . 178 178 . 7.648   39.320 -4.701  1.00 50.00 .
ATOM   1332 CE  CE  . MET MET . . 178 178 . 8.694   40.566 -5.419  1.00 50.00 .
ATOM   1333 N   N   . GLU GLU . . 179 179 . 10.309  37.212 -8.746  1.00 50.00 .
ATOM   1334 CA  CA  . GLU GLU . . 179 179 . 11.738  36.949 -8.531  1.00 50.00 .
ATOM   1335 C   C   . GLU GLU . . 179 179 . 12.277  37.843 -7.413  1.00 50.00 .
ATOM   1336 O   O   . GLU GLU . . 179 179 . 12.454  39.048 -7.588  1.00 50.00 .
ATOM   1337 CB  CB  . GLU GLU . . 179 179 . 12.540  37.097 -9.829  1.00 50.00 .
ATOM   1338 CG  CG  . GLU GLU . . 179 179 . 14.046  36.954 -9.573  1.00 50.00 .
ATOM   1339 CD  CD  . GLU GLU . . 179 179 . 14.709  36.006 -10.564 1.00 50.00 .
ATOM   1340 OE1 OE1 . GLU GLU . . 179 179 . 14.992  36.476 -11.686 1.00 50.00 .
ATOM   1341 OE2 OE2 . GLU GLU . . 179 179 . 14.881  34.828 -10.183 1.00 50.00 .
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ATOM   1342 N   N   . GLY GLY . . 180 180 . 12.509  37.176 -6.280  1.00 50.00 .
ATOM   1343 CA  CA  . GLY GLY . . 180 180 . 13.228  37.779 -5.153  1.00 50.00 .
ATOM   1344 C   C   . GLY GLY . . 180 180 . 14.616  38.215 -5.601  1.00 50.00 .
ATOM   1345 O   O   . GLY GLY . . 180 180 . 14.859  39.388 -5.874  1.00 50.00 .
ATOM   1346 N   N   . ILE ILE . . 181 181 . 15.452  37.189 -5.730  1.00 50.00 .
ATOM   1347 CA  CA  . ILE ILE . . 181 181 . 16.855  37.306 -6.194  1.00 50.00 .
ATOM   1348 C   C   . ILE ILE . . 181 181 . 17.297  35.968 -6.841  1.00 50.00 .
ATOM   1349 O   O   . ILE ILE . . 181 181 . 17.943  35.158 -6.138  1.00 50.00 .
ATOM   1350 CB  CB  . ILE ILE . . 181 181 . 17.804  37.728 -5.040  1.00 50.00 .
ATOM   1351 CG1 CG1 . ILE ILE . . 181 181 . 17.355  39.009 -4.310  1.00 50.00 .
ATOM   1352 CG2 CG2 . ILE ILE . . 181 181 . 19.221  37.946 -5.597  1.00 50.00 .
ATOM   1353 CD1 CD1 . ILE ILE . . 181 181 . 18.099  39.349 -3.013  1.00 50.00 .
ATOM   1354 OXT OXT . ILE ILE . . 181 181 . 16.896  35.748 -8.005  1.00 99.99 .
ATOM   1355 N   N   . GLY GLY A A -1  -1  . 4.352   39.773 -22.016 1.00 70.16 .
ATOM   1356 CA  CA  . GLY GLY A A -1  -1  . 4.332   38.478 -21.278 1.00 69.23 .
ATOM   1357 C   C   . GLY GLY A A -1  -1  . 4.803   37.322 -22.139 1.00 71.79 .
ATOM   1358 O   O   . GLY GLY A A -1  -1  . 4.031   36.415 -22.462 1.00 73.68 .
ATOM   1359 N   N   . SER SER A A 0   0   . 6.076   37.355 -22.519 1.00 68.23 .
ATOM   1360 CA  CA  . SER SER A A 0   0   . 6.649   36.302 -23.346 1.00 64.14 .
ATOM   1361 C   C   . SER SER A A 0   0   . 6.138   34.932 -22.906 1.00 67.09 .
ATOM   1362 O   O   . SER SER A A 0   0   . 6.129   34.610 -21.715 1.00 67.64 .
ATOM   1363 CB  CB  . SER SER A A 0   0   . 8.173   36.347 -23.261 1.00 50.31 .
ATOM   1364 OG  OG  . SER SER A A 0   0   . 8.749   35.245 -23.955 1.00 50.92 .
ATOM   1365 N   N   . ASP ASP A A 22  22  . 5.712   34.130 -23.876 1.00 68.26 .
ATOM   1366 CA  CA  . ASP ASP A A 22  22  . 5.174   32.800 -23.602 1.00 66.93 .
ATOM   1367 C   C   . ASP ASP A A 22  22  . 6.250   31.830 -23.126 1.00 63.29 .
ATOM   1368 O   O   . ASP ASP A A 22  22  . 6.002   30.630 -23.015 1.00 63.81 .
ATOM   1369 CB  CB  . ASP ASP A A 22  22  . 4.504   32.239 -24.862 1.00 71.28 .
ATOM   1370 CG  CG  . ASP ASP A A 22  22  . 3.741   33.300 -25.639 1.00 75.97 .
ATOM   1371 OD1 OD1 . ASP ASP A A 22  22  . 2.933   34.028 -25.021 1.00 80.52 .
ATOM   1372 OD2 OD2 . ASP ASP A A 22  22  . 3.947   33.403 -26.870 1.00 70.15 .
ATOM   1373 N   N   . GLU GLU A A 23  23  . 7.440   32.350 -22.844 1.00 59.67 .
ATOM   1374 CA  CA  . GLU GLU A A 23  23  . 8.542   31.514 -22.390 1.00 59.63 .
ATOM   1375 C   C   . GLU GLU A A 23  23  . 9.183   32.069 -21.118 1.00 56.26 .
ATOM   1376 O   O   . GLU GLU A A 23  23  . 10.366  31.858 -20.858 1.00 53.06 .
ATOM   1377 CB  CB  . GLU GLU A A 23  23  . 9.586   31.372 -23.510 1.00 65.14 .
ATOM   1378 CG  CG  . GLU GLU A A 23  23  . 10.675  30.342 -23.232 1.00 73.21 .
ATOM   1379 CD  CD  . GLU GLU A A 23  23  . 11.429  29.922 -24.483 1.00 82.25 .
ATOM   1380 OE1 OE1 . GLU GLU A A 23  23  . 12.382  29.120 -24.363 1.00 86.14 .
ATOM   1381 OE2 OE2 . GLU GLU A A 23  23  . 11.070  30.386 -25.587 1.00 85.38 .
ATOM   1382 N   N   . ASP ASP A A 24  24  . 8.388   32.783 -20.327 1.00 55.49 .
ATOM   1383 CA  CA  . ASP ASP A A 24  24  . 8.862   33.345 -19.066 1.00 55.29 .
ATOM   1384 C   C   . ASP ASP A A 24  24  . 8.437   32.382 -17.960 1.00 51.70 .
ATOM   1385 O   O   . ASP ASP A A 24  24  . 8.057   32.795 -16.863 1.00 53.18 .
ATOM   1386 CB  CB  . ASP ASP A A 24  24  . 8.244   34.731 -18.825 1.00 53.20 .
ATOM   1387 CG  CG  . ASP ASP A A 24  24  . 6.734   34.681 -18.600 1.00 63.37 .
ATOM   1388 OD1 OD1 . ASP ASP A A 24  24  . 6.129   35.757 -18.400 1.00 67.71 .
ATOM   1389 OD2 OD2 . ASP ASP A A 24  24  . 6.149   33.576 -18.621 1.00 66.79 .
ATOM   1390 N   N   . PHE PHE A A 25  25  . 8.509   31.090 -18.262 1.00 46.37 .
ATOM   1391 CA  CA  . PHE PHE A A 25  25  . 8.104   30.054 -17.322 1.00 43.02 .
ATOM   1392 C   C   . PHE PHE A A 25  25  . 8.962   29.901 -16.076 1.00 37.35 .
ATOM   1393 O   O   . PHE PHE A A 25  25  . 10.134  30.264 -16.053 1.00 34.94 .
ATOM   1394 CB  CB  . PHE PHE A A 25  25  . 8.004   28.704 -18.039 1.00 48.23 .
ATOM   1395 CG  CG  . PHE PHE A A 25  25  . 9.180   28.392 -18.916 1.00 50.75 .
ATOM   1396 CD1 CD1 . PHE PHE A A 25  25  . 10.451  28.238 -18.376 1.00 54.74 .
ATOM   1397 CD2 CD2 . PHE PHE A A 25  25  . 9.014   28.247 -20.288 1.00 55.97 .
ATOM   1398 CE1 CE1 . PHE PHE A A 25  25  . 11.546  27.942 -19.195 1.00 64.32 .
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ATOM   1399 CE2 CE2 . PHE PHE A A 25  25  . 10.099  27.952 -21.113 1.00 60.96 .
ATOM   1400 CZ  CZ  . PHE PHE A A 25  25  . 11.368  27.799 -20.564 1.00 60.70 .
ATOM   1401 N   N   . VAL VAL A A 26  26  . 8.342   29.357 -15.034 1.00 35.80 .
ATOM   1402 CA  CA  . VAL VAL A A 26  26  . 9.001   29.114 -13.760 1.00 27.22 .
ATOM   1403 C   C   . VAL VAL A A 26  26  . 9.806   27.837 -13.934 1.00 30.24 .
ATOM   1404 O   O   . VAL VAL A A 26  26  . 9.303   26.855 -14.481 1.00 28.59 .
ATOM   1405 CB  CB  . VAL VAL A A 26  26  . 7.961   28.920 -12.643 1.00 29.39 .
ATOM   1406 CG1 CG1 . VAL VAL A A 26  26  . 8.657   28.697 -11.322 1.00 24.16 .
ATOM   1407 CG2 CG2 . VAL VAL A A 26  26  . 7.047   30.136 -12.573 1.00 28.87 .
ATOM   1408 N   N   . THR THR A A 27  27  . 11.048  27.845 -13.462 1.00 28.75 .
ATOM   1409 CA  CA  . THR THR A A 27  27  . 11.924  26.692 -13.627 1.00 27.43 .
ATOM   1410 C   C   . THR THR A A 27  27  . 12.291  25.947 -12.351 1.00 24.58 .
ATOM   1411 O   O   . THR THR A A 27  27  . 12.247  26.494 -11.248 1.00 0.00  .
ATOM   1412 CB  CB  . THR THR A A 27  27  . 13.221  27.110 -14.340 1.00 29.95 .
ATOM   1413 OG1 OG1 . THR THR A A 27  27  . 13.958  28.006 -13.500 1.00 34.53 .
ATOM   1414 CG2 CG2 . THR THR A A 27  27  . 12.897  27.816 -15.655 1.00 30.45 .
ATOM   1415 N   N   . CYS CYS A A 28  28  . 12.678  24.691 -12.520 1.00 18.60 .
ATOM   1416 CA  CA  . CYS CYS A A 28  28  . 13.041  23.850 -11.395 1.00 21.71 .
ATOM   1417 C   C   . CYS CYS A A 28  28  . 14.303  24.370 -10.707 1.00 0.00  .
ATOM   1418 O   O   . CYS CYS A A 28  28  . 15.269  24.735 -11.365 1.00 23.94 .
ATOM   1419 CB  CB  . CYS CYS A A 28  28  . 13.237  22.421 -11.884 1.00 25.22 .
ATOM   1420 SG  SG  . CYS CYS A A 28  28  . 12.735  21.174 -10.668 1.00 34.40 .
ATOM   1421 N   N   . TYR TYR A A 29  29  . 14.266  24.401 -9.379  1.00 18.10 .
ATOM   1422 CA  CA  . TYR TYR A A 29  29  . 15.354  24.896 -8.536  1.00 22.23 .
ATOM   1423 C   C   . TYR TYR A A 29  29  . 15.496  26.413 -8.554  1.00 23.58 .
ATOM   1424 O   O   . TYR TYR A A 29  29  . 16.460  26.960 -8.018  1.00 22.74 .
ATOM   1425 CB  CB  . TYR TYR A A 29  29  . 16.689  24.237 -8.902  1.00 24.02 .
ATOM   1426 CG  CG  . TYR TYR A A 29  29  . 16.721  22.772 -8.547  1.00 19.27 .
ATOM   1427 CD1 CD1 . TYR TYR A A 29  29  . 16.456  22.348 -7.246  1.00 19.92 .
ATOM   1428 CD2 CD2 . TYR TYR A A 29  29  . 16.975  21.808 -9.516  1.00 21.97 .
ATOM   1429 CE1 CE1 . TYR TYR A A 29  29  . 16.438  20.996 -6.922  1.00 29.97 .
ATOM   1430 CE2 CE2 . TYR TYR A A 29  29  . 16.960  20.459 -9.206  1.00 27.20 .
ATOM   1431 CZ  CZ  . TYR TYR A A 29  29  . 16.692  20.057 -7.910  1.00 27.70 .
ATOM   1432 OH  OH  . TYR TYR A A 29  29  . 16.678  18.717 -7.612  1.00 36.46 .
ATOM   1433 N   N   . SER SER A A 30  30  . 14.547  27.100 -9.178  1.00 23.89 .
ATOM   1434 CA  CA  . SER SER A A 30  30  . 14.603  28.554 -9.184  1.00 20.88 .
ATOM   1435 C   C   . SER SER A A 30  30  . 14.223  28.977 -7.762  1.00 26.23 .
ATOM   1436 O   O   . SER SER A A 30  30  . 13.589  28.212 -7.034  1.00 20.15 .
ATOM   1437 CB  CB  . SER SER A A 30  30  . 13.612  29.125 -10.204 1.00 32.39 .
ATOM   1438 OG  OG  . SER SER A A 30  30  . 12.296  28.675 -9.943  1.00 34.78 .
ATOM   1439 N   N   . VAL VAL A A 31  31  . 14.632  30.175 -7.357  1.00 20.82 .
ATOM   1440 CA  CA  . VAL VAL A A 31  31  . 14.320  30.683 -6.024  1.00 20.03 .
ATOM   1441 C   C   . VAL VAL A A 31  31  . 13.468  31.920 -6.213  1.00 22.76 .
ATOM   1442 O   O   . VAL VAL A A 31  31  . 13.848  32.824 -6.956  1.00 19.69 .
ATOM   1443 CB  CB  . VAL VAL A A 31  31  . 15.586  31.095 -5.248  1.00 18.30 .
ATOM   1444 CG1 CG1 . VAL VAL A A 31  31  . 15.198  31.706 -3.917  1.00 16.03 .
ATOM   1445 CG2 CG2 . VAL VAL A A 31  31  . 16.472  29.894 -5.021  1.00 17.26 .
ATOM   1446 N   N   . LEU LEU A A 32  32  . 12.329  31.973 -5.538  1.00 19.71 .
ATOM   1447 CA  CA  . LEU LEU A A 32  32  . 11.445  33.118 -5.691  1.00 25.12 .
ATOM   1448 C   C   . LEU LEU A A 32  32  . 10.511  33.343 -4.511  1.00 18.61 .
ATOM   1449 O   O   . LEU LEU A A 32  32  . 10.534  32.593 -3.537  1.00 17.80 .
ATOM   1450 CB  CB  . LEU LEU A A 32  32  . 10.644  32.946 -6.984  1.00 38.48 .
ATOM   1451 CG  CG  . LEU LEU A A 32  32  . 10.229  31.502 -7.265  1.00 38.83 .
ATOM   1452 CD1 CD1 . LEU LEU A A 32  32  . 9.129   31.120 -6.319  1.00 39.95 .
ATOM   1453 CD2 CD2 . LEU LEU A A 32  32  . 9.775   31.336 -8.701  1.00 44.74 .
ATOM   1454 N   N   . LYS LYS A A 33  33  . 9.708   34.398 -4.606  1.00 14.81 .
ATOM   1455 CA  CA  . LYS LYS A A 33  33  . 8.741   34.747 -3.574  1.00 18.14 .
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ATOM   1456 C   C   . LYS LYS A A 33  33  . 7.343   34.433 -4.086  1.00 16.97 .
ATOM   1457 O   O   . LYS LYS A A 33  33  . 7.067   34.584 -5.277  1.00 20.25 .
ATOM   1458 CB  CB  . LYS LYS A A 33  33  . 8.816   36.242 -3.251  1.00 17.22 .
ATOM   1459 CG  CG  . LYS LYS A A 33  33  . 10.159  36.721 -2.744  1.00 26.26 .
ATOM   1460 CD  CD  . LYS LYS A A 33  33  . 10.261  38.234 -2.901  1.00 33.27 .
ATOM   1461 CE  CE  . LYS LYS A A 33  33  . 11.630  38.746 -2.511  1.00 51.37 .
ATOM   1462 NZ  NZ  . LYS LYS A A 33  33  . 11.894  40.092 -3.111  1.00 62.26 .
ATOM   1463 N   N   . PHE PHE A A 34  34  . 6.470   33.991 -3.186  1.00 18.51 .
ATOM   1464 CA  CA  . PHE PHE A A 34  34  . 5.088   33.683 -3.525  1.00 17.35 .
ATOM   1465 C   C   . PHE PHE A A 34  34  . 4.221   34.766 -2.899  1.00 25.06 .
ATOM   1466 O   O   . PHE PHE A A 34  34  . 4.102   34.849 -1.676  1.00 20.95 .
ATOM   1467 CB  CB  . PHE PHE A A 34  34  . 4.684   32.308 -2.983  1.00 16.74 .
ATOM   1468 CG  CG  . PHE PHE A A 34  34  . 5.322   31.158 -3.719  1.00 22.57 .
ATOM   1469 CD1 CD1 . PHE PHE A A 34  34  . 6.654   30.824 -3.498  1.00 31.37 .
ATOM   1470 CD2 CD2 . PHE PHE A A 34  34  . 4.606   30.449 -4.675  1.00 30.93 .
ATOM   1471 CE1 CE1 . PHE PHE A A 34  34  . 7.262   29.807 -4.223  1.00 29.22 .
ATOM   1472 CE2 CE2 . PHE PHE A A 34  34  . 5.204   29.429 -5.410  1.00 27.96 .
ATOM   1473 CZ  CZ  . PHE PHE A A 34  34  . 6.534   29.109 -5.185  1.00 27.78 .
ATOM   1474 N   N   . ILE ILE A A 35  35  . 3.615   35.593 -3.744  1.00 20.74 .
ATOM   1475 CA  CA  . ILE ILE A A 35  35  . 2.781   36.694 -3.273  1.00 21.08 .
ATOM   1476 C   C   . ILE ILE A A 35  35  . 1.287   36.378 -3.371  1.00 21.63 .
ATOM   1477 O   O   . ILE ILE A A 35  35  . 0.807   35.922 -4.411  1.00 20.71 .
ATOM   1478 CB  CB  . ILE ILE A A 35  35  . 3.068   37.973 -4.093  1.00 16.30 .
ATOM   1479 CG1 CG1 . ILE ILE A A 35  35  . 4.552   38.337 -3.974  1.00 25.64 .
ATOM   1480 CG2 CG2 . ILE ILE A A 35  35  . 2.178   39.126 -3.611  1.00 19.77 .
ATOM   1481 CD1 CD1 . ILE ILE A A 35  35  . 4.932   39.556 -4.772  1.00 41.55 .
ATOM   1482 N   N   . ASN ASN A A 36  36  . 0.546   36.620 -2.297  1.00 16.26 .
ATOM   1483 CA  CA  . ASN ASN A A 36  36  . -0.883  36.351 -2.354  1.00 16.95 .
ATOM   1484 C   C   . ASN ASN A A 36  36  . -1.545  37.436 -3.193  1.00 19.05 .
ATOM   1485 O   O   . ASN ASN A A 36  36  . -1.487  38.626 -2.868  1.00 21.03 .
ATOM   1486 CB  CB  . ASN ASN A A 36  36  . -1.483  36.291 -0.951  1.00 16.89 .
ATOM   1487 CG  CG  . ASN ASN A A 36  36  . -2.917  35.806 -0.964  1.00 14.38 .
ATOM   1488 OD1 OD1 . ASN ASN A A 36  36  . -3.844  36.593 -1.107  1.00 18.19 .
ATOM   1489 ND2 ND2 . ASN ASN A A 36  36  . -3.103  34.501 -0.829  1.00 19.52 .
ATOM   1490 N   N   . ALA ALA A A 37  37  . -2.153  37.010 -4.294  1.00 29.08 .
ATOM   1491 CA  CA  . ALA ALA A A 37  37  . -2.807  37.910 -5.240  1.00 22.33 .
ATOM   1492 C   C   . ALA ALA A A 37  37  . -3.721  38.987 -4.664  1.00 27.07 .
ATOM   1493 O   O   . ALA ALA A A 37  37  . -3.595  40.164 -5.002  1.00 28.70 .
ATOM   1494 CB  CB  . ALA ALA A A 37  37  . -3.582  37.090 -6.264  1.00 26.81 .
ATOM   1495 N   N   . ASN ASN A A 38  38  . -4.643  38.597 -3.796  1.00 27.01 .
ATOM   1496 CA  CA  . ASN ASN A A 38  38  . -5.590  39.550 -3.241  1.00 29.10 .
ATOM   1497 C   C   . ASN ASN A A 38  38  . -5.125  40.351 -2.011  1.00 32.55 .
ATOM   1498 O   O   . ASN ASN A A 38  38  . -5.605  41.466 -1.771  1.00 22.14 .
ATOM   1499 CB  CB  . ASN ASN A A 38  38  . -6.902  38.814 -2.954  1.00 33.75 .
ATOM   1500 CG  CG  . ASN ASN A A 38  38  . -7.560  38.263 -4.231  1.00 39.89 .
ATOM   1501 OD1 OD1 . ASN ASN A A 38  38  . -8.464  37.423 -4.168  1.00 39.69 .
ATOM   1502 ND2 ND2 . ASN ASN A A 38  38  . -7.114  38.748 -5.390  1.00 27.62 .
ATOM   1503 N   N   . ASP ASP A A 39  39  . -4.183  39.804 -1.247  1.00 29.19 .
ATOM   1504 CA  CA  . ASP ASP A A 39  39  . -3.686  40.494 -0.056  1.00 27.69 .
ATOM   1505 C   C   . ASP ASP A A 39  39  . -2.323  41.152 -0.267  1.00 22.47 .
ATOM   1506 O   O   . ASP ASP A A 39  39  . -1.992  42.132 0.401   1.00 25.32 .
ATOM   1507 CB  CB  . ASP ASP A A 39  39  . -3.600  39.518 1.124   1.00 33.35 .
ATOM   1508 CG  CG  . ASP ASP A A 39  39  . -3.242  40.209 2.419   1.00 34.85 .
ATOM   1509 OD1 OD1 . ASP ASP A A 39  39  . -3.853  41.260 2.708   1.00 38.48 .
ATOM   1510 OD2 OD2 . ASP ASP A A 39  39  . -2.365  39.706 3.149   1.00 21.75 .
ATOM   1511 N   N   . GLY GLY A A 40  40  . -1.527  40.609 -1.182  1.00 18.58 .
ATOM   1512 CA  CA  . GLY GLY A A 40  40  . -0.209  41.178 -1.440  1.00 20.86 .
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ATOM   1513 C   C   . GLY GLY A A 40  40  . 0.899   40.694 -0.506  1.00 20.31 .
ATOM   1514 O   O   . GLY GLY A A 40  40  . 2.064   41.045 -0.690  1.00 23.40 .
ATOM   1515 N   N   . SER SER A A 41  41  . 0.546   39.888 0.493   1.00 17.03 .
ATOM   1516 CA  CA  . SER SER A A 41  41  . 1.526   39.361 1.444   1.00 18.80 .
ATOM   1517 C   C   . SER SER A A 41  41  . 2.366   38.237 0.821   1.00 21.39 .
ATOM   1518 O   O   . SER SER A A 41  41  . 1.910   37.549 -0.095  1.00 18.73 .
ATOM   1519 CB  CB  . SER SER A A 41  41  . 0.805   38.843 2.689   1.00 12.61 .
ATOM   1520 OG  OG  . SER SER A A 41  41  . -0.199  37.908 2.341   1.00 20.94 .
ATOM   1521 N   N   . ARG ARG A A 42  42  . 3.587   38.053 1.322   1.00 15.64 .
ATOM   1522 CA  CA  . ARG ARG A A 42  42  . 4.486   37.028 0.795   1.00 16.17 .
ATOM   1523 C   C   . ARG ARG A A 42  42  . 4.558   35.794 1.676   1.00 17.24 .
ATOM   1524 O   O   . ARG ARG A A 42  42  . 4.645   35.911 2.901   1.00 14.95 .
ATOM   1525 CB  CB  . ARG ARG A A 42  42  . 5.903   37.581 0.632   1.00 18.32 .
ATOM   1526 CG  CG  . ARG ARG A A 42  42  . 6.061   38.608 -0.464  1.00 19.44 .
ATOM   1527 CD  CD  . ARG ARG A A 42  42  . 5.686   39.995 0.012   1.00 31.46 .
ATOM   1528 NE  NE  . ARG ARG A A 42  42  . 5.611   40.944 -1.096  1.00 43.52 .
ATOM   1529 CZ  CZ  . ARG ARG A A 42  42  . 6.634   41.277 -1.873  1.00 45.62 .
ATOM   1530 NH1 NH1 . ARG ARG A A 42  42  . 7.833   40.743 -1.671  1.00 53.39 .
ATOM   1531 NH2 NH2 . ARG ARG A A 42  42  . 6.454   42.146 -2.858  1.00 0.00  .
ATOM   1532 N   N   . LEU LEU A A 43  43  . 4.529   34.617 1.050   1.00 11.44 .
ATOM   1533 CA  CA  . LEU LEU A A 43  43  . 4.613   33.352 1.791   1.00 10.82 .
ATOM   1534 C   C   . LEU LEU A A 43  43  . 5.896   33.452 2.592   1.00 12.80 .
ATOM   1535 O   O   . LEU LEU A A 43  43  . 6.936   33.881 2.067   1.00 11.97 .
ATOM   1536 CB  CB  . LEU LEU A A 43  43  . 4.653   32.159 0.837   1.00 12.78 .
ATOM   1537 CG  CG  . LEU LEU A A 43  43  . 4.489   30.796 1.521   1.00 10.91 .
ATOM   1538 CD1 CD1 . LEU LEU A A 43  43  . 3.177   30.749 2.269   1.00 13.86 .
ATOM   1539 CD2 CD2 . LEU LEU A A 43  43  . 4.543   29.692 0.480   1.00 9.77  .
ATOM   1540 N   N   . HIS HIS A A 44  44  . 5.850   33.008 3.843   1.00 13.24 .
ATOM   1541 CA  CA  . HIS HIS A A 44  44  . 6.996   33.218 4.718   1.00 10.40 .
ATOM   1542 C   C   . HIS HIS A A 44  44  . 7.148   32.153 5.801   1.00 17.84 .
ATOM   1543 O   O   . HIS HIS A A 44  44  . 6.184   31.470 6.173   1.00 15.55 .
ATOM   1544 CB  CB  . HIS HIS A A 44  44  . 6.769   34.605 5.348   1.00 15.13 .
ATOM   1545 CG  CG  . HIS HIS A A 44  44  . 7.875   35.107 6.228   1.00 22.74 .
ATOM   1546 ND1 ND1 . HIS HIS A A 44  44  . 7.931   34.839 7.578   1.00 28.39 .
ATOM   1547 CD2 CD2 . HIS HIS A A 44  44  . 8.902   35.953 5.970   1.00 16.33 .
ATOM   1548 CE1 CE1 . HIS HIS A A 44  44  . 8.942   35.500 8.116   1.00 20.73 .
ATOM   1549 NE2 NE2 . HIS HIS A A 44  44  . 9.546   36.184 7.161   1.00 23.95 .
ATOM   1550 N   N   . SER SER A A 45  45  . 8.370   31.994 6.286   1.00 16.17 .
ATOM   1551 CA  CA  . SER SER A A 45  45  . 8.607   31.074 7.384   1.00 16.77 .
ATOM   1552 C   C   . SER SER A A 45  45  . 9.720   31.632 8.254   1.00 24.09 .
ATOM   1553 O   O   . SER SER A A 45  45  . 10.345  32.633 7.904   1.00 20.56 .
ATOM   1554 CB  CB  . SER SER A A 45  45  . 8.947   29.677 6.901   1.00 15.48 .
ATOM   1555 OG  OG  . SER SER A A 45  45  . 8.748   28.766 7.974   1.00 19.35 .
ATOM   1556 N   N   . HIS HIS A A 46  46  . 9.977   30.976 9.379   1.00 20.47 .
ATOM   1557 CA  CA  . HIS HIS A A 46  46  . 10.971  31.469 10.321  1.00 24.37 .
ATOM   1558 C   C   . HIS HIS A A 46  46  . 11.146  30.434 11.421  1.00 24.45 .
ATOM   1559 O   O   . HIS HIS A A 46  46  . 10.519  29.384 11.381  1.00 22.36 .
ATOM   1560 CB  CB  . HIS HIS A A 46  46  . 10.445  32.778 10.893  1.00 21.36 .
ATOM   1561 CG  CG  . HIS HIS A A 46  46  . 8.963   32.768 11.106  1.00 18.89 .
ATOM   1562 ND1 ND1 . HIS HIS A A 46  46  . 8.206   33.917 11.166  1.00 21.40 .
ATOM   1563 CD2 CD2 . HIS HIS A A 46  46  . 8.096   31.739 11.273  1.00 20.77 .
ATOM   1564 CE1 CE1 . HIS HIS A A 46  46  . 6.938   33.598 11.362  1.00 18.99 .
ATOM   1565 NE2 NE2 . HIS HIS A A 46  46  . 6.845   32.282 11.431  1.00 16.44 .
ATOM   1566 N   N   . ASP ASP A A 47  47  . 11.984  30.734 12.407  1.00 29.95 .
ATOM   1567 CA  CA  . ASP ASP A A 47  47  . 12.236  29.795 13.496  1.00 30.32 .
ATOM   1568 C   C   . ASP ASP A A 47  47  . 11.204  29.923 14.609  1.00 34.98 .
ATOM   1569 O   O   . ASP ASP A A 47  47  . 11.551  29.952 15.793  1.00 47.26 .
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ATOM   1570 CB  CB  . ASP ASP A A 47  47  . 13.638  30.019 14.077  1.00 34.20 .
ATOM   1571 CG  CG  . ASP ASP A A 47  47  . 14.169  28.795 14.804  1.00 47.14 .
ATOM   1572 OD1 OD1 . ASP ASP A A 47  47  . 13.375  28.120 15.494  1.00 52.03 .
ATOM   1573 OD2 OD2 . ASP ASP A A 47  47  . 15.382  28.511 14.691  1.00 42.39 .
ATOM   1574 N   N   . VAL VAL A A 48  48  . 9.936   30.016 14.229  1.00 32.28 .
ATOM   1575 CA  CA  . VAL VAL A A 48  48  . 8.846   30.122 15.191  1.00 24.97 .
ATOM   1576 C   C   . VAL VAL A A 48  48  . 7.950   28.917 14.959  1.00 36.24 .
ATOM   1577 O   O   . VAL VAL A A 48  48  . 7.397   28.756 13.873  1.00 30.64 .
ATOM   1578 CB  CB  . VAL VAL A A 48  48  . 8.021   31.412 14.975  1.00 29.49 .
ATOM   1579 CG1 CG1 . VAL VAL A A 48  48  . 6.831   31.438 15.922  1.00 22.95 .
ATOM   1580 CG2 CG2 . VAL VAL A A 48  48  . 8.901   32.641 15.196  1.00 27.27 .
ATOM   1581 N   N   . LYS LYS A A 49  49  . 7.810   28.072 15.978  1.00 36.26 .
ATOM   1582 CA  CA  . LYS LYS A A 49  49  . 6.994   26.871 15.862  1.00 33.71 .
ATOM   1583 C   C   . LYS LYS A A 49  49  . 5.566   27.049 16.356  1.00 36.36 .
ATOM   1584 O   O   . LYS LYS A A 49  49  . 5.283   27.932 17.164  1.00 34.90 .
ATOM   1585 CB  CB  . LYS LYS A A 49  49  . 7.654   25.713 16.620  1.00 39.65 .
ATOM   1586 CG  CG  . LYS LYS A A 49  49  . 9.061   25.390 16.128  1.00 51.06 .
ATOM   1587 CD  CD  . LYS LYS A A 49  49  . 9.519   24.001 16.545  1.00 50.82 .
ATOM   1588 CE  CE  . LYS LYS A A 49  49  . 10.858  23.657 15.902  1.00 49.46 .
ATOM   1589 NZ  NZ  . LYS LYS A A 49  49  . 11.276  22.243 16.151  1.00 54.24 .
ATOM   1590 N   N   . TYR TYR A A 50  50  . 4.667   26.208 15.851  1.00 33.34 .
ATOM   1591 CA  CA  . TYR TYR A A 50  50  . 3.268   26.248 16.249  1.00 33.99 .
ATOM   1592 C   C   . TYR TYR A A 50  50  . 3.149   25.726 17.670  1.00 41.58 .
ATOM   1593 O   O   . TYR TYR A A 50  50  . 3.989   24.957 18.136  1.00 42.00 .
ATOM   1594 CB  CB  . TYR TYR A A 50  50  . 2.406   25.359 15.348  1.00 38.38 .
ATOM   1595 CG  CG  . TYR TYR A A 50  50  . 2.135   25.894 13.958  1.00 37.31 .
ATOM   1596 CD1 CD1 . TYR TYR A A 50  50  . 1.508   27.128 13.771  1.00 34.21 .
ATOM   1597 CD2 CD2 . TYR TYR A A 50  50  . 2.462   25.141 12.829  1.00 30.81 .
ATOM   1598 CE1 CE1 . TYR TYR A A 50  50  . 1.209   27.598 12.489  1.00 33.56 .
ATOM   1599 CE2 CE2 . TYR TYR A A 50  50  . 2.169   25.600 11.546  1.00 29.03 .
ATOM   1600 CZ  CZ  . TYR TYR A A 50  50  . 1.543   26.827 11.382  1.00 34.74 .
ATOM   1601 OH  OH  . TYR TYR A A 50  50  . 1.243   27.278 10.116  1.00 27.67 .
ATOM   1602 N   N   . GLY GLY A A 51  51  . 2.093   26.149 18.349  1.00 45.04 .
ATOM   1603 CA  CA  . GLY GLY A A 51  51  . 1.847   25.708 19.704  1.00 52.07 .
ATOM   1604 C   C   . GLY GLY A A 51  51  . 0.415   25.227 19.728  1.00 58.90 .
ATOM   1605 O   O   . GLY GLY A A 51  51  . -0.384  25.672 20.551  1.00 62.54 .
ATOM   1606 N   N   . SER SER A A 52  52  . 0.088   24.322 18.809  1.00 59.77 .
ATOM   1607 CA  CA  . SER SER A A 52  52  . -1.265  23.794 18.716  1.00 60.84 .
ATOM   1608 C   C   . SER SER A A 52  52  . -1.370  22.455 17.980  1.00 60.63 .
ATOM   1609 O   O   . SER SER A A 52  52  . -1.536  21.410 18.608  1.00 69.40 .
ATOM   1610 CB  CB  . SER SER A A 52  52  . -2.173  24.827 18.042  1.00 62.63 .
ATOM   1611 OG  OG  . SER SER A A 52  52  . -1.625  25.266 16.806  1.00 61.55 .
ATOM   1612 N   N   . GLY GLY A A 53  53  . -1.275  22.488 16.655  1.00 53.19 .
ATOM   1613 CA  CA  . GLY GLY A A 53  53  . -1.394  21.270 15.869  1.00 48.98 .
ATOM   1614 C   C   . GLY GLY A A 53  53  . -0.265  20.262 15.992  1.00 51.02 .
ATOM   1615 O   O   . GLY GLY A A 53  53  . -0.210  19.481 16.942  1.00 46.29 .
ATOM   1616 N   N   . SER SER A A 54  54  . 0.634   20.270 15.014  1.00 43.61 .
ATOM   1617 CA  CA  . SER SER A A 54  54  . 1.767   19.351 14.996  1.00 38.34 .
ATOM   1618 C   C   . SER SER A A 54  54  . 2.923   19.916 15.804  1.00 38.29 .
ATOM   1619 O   O   . SER SER A A 54  54  . 3.861   19.200 16.149  1.00 39.48 .
ATOM   1620 CB  CB  . SER SER A A 54  54  . 2.235   19.122 13.556  1.00 45.61 .
ATOM   1621 OG  OG  . SER SER A A 54  54  . 2.833   20.298 13.025  1.00 35.65 .
ATOM   1622 N   N   . GLY GLY A A 55  55  . 2.852   21.208 16.099  1.00 36.84 .
ATOM   1623 CA  CA  . GLY GLY A A 55  55  . 3.919   21.840 16.846  1.00 34.17 .
ATOM   1624 C   C   . GLY GLY A A 55  55  . 5.156   22.009 15.982  1.00 31.24 .
ATOM   1625 O   O   . GLY GLY A A 55  55  . 6.256   22.214 16.495  1.00 27.38 .
ATOM   1626 N   N   . GLN GLN A A 56  56  . 4.985   21.910 14.666  1.00 27.92 .
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ATOM   1627 CA  CA  . GLN GLN A A 56  56  . 6.106   22.073 13.742  1.00 27.72 .
ATOM   1628 C   C   . GLN GLN A A 56  56  . 6.299   23.558 13.429  1.00 23.04 .
ATOM   1629 O   O   . GLN GLN A A 56  56  . 5.508   24.392 13.856  1.00 22.32 .
ATOM   1630 CB  CB  . GLN GLN A A 56  56  . 5.847   21.303 12.445  1.00 25.38 .
ATOM   1631 CG  CG  . GLN GLN A A 56  56  . 5.633   19.817 12.645  1.00 37.82 .
ATOM   1632 CD  CD  . GLN GLN A A 56  56  . 5.376   19.088 11.342  1.00 39.23 .
ATOM   1633 OE1 OE1 . GLN GLN A A 56  56  . 4.722   19.612 10.446  1.00 30.71 .
ATOM   1634 NE2 NE2 . GLN GLN A A 56  56  . 5.877   17.862 11.239  1.00 40.25 .
ATOM   1635 N   N   . GLN GLN A A 57  57  . 7.346   23.875 12.676  1.00 23.53 .
ATOM   1636 CA  CA  . GLN GLN A A 57  57  . 7.651   25.256 12.309  1.00 23.22 .
ATOM   1637 C   C   . GLN GLN A A 57  57  . 6.471   25.867 11.568  1.00 22.82 .
ATOM   1638 O   O   . GLN GLN A A 57  57  . 5.940   25.264 10.640  1.00 23.62 .
ATOM   1639 CB  CB  . GLN GLN A A 57  57  . 8.891   25.295 11.421  1.00 21.79 .
ATOM   1640 CG  CG  . GLN GLN A A 57  57  . 9.487   26.673 11.250  1.00 18.27 .
ATOM   1641 CD  CD  . GLN GLN A A 57  57  . 10.762  26.637 10.446  1.00 29.61 .
ATOM   1642 OE1 OE1 . GLN GLN A A 57  57  . 10.787  27.014 9.267   1.00 25.89 .
ATOM   1643 NE2 NE2 . GLN GLN A A 57  57  . 11.832  26.162 11.068  1.00 22.57 .
ATOM   1644 N   N   . SER SER A A 58  58  . 6.068   27.065 11.978  1.00 24.36 .
ATOM   1645 CA  CA  . SER SER A A 58  58  . 4.935   27.735 11.357  1.00 20.75 .
ATOM   1646 C   C   . SER SER A A 58  58  . 5.238   28.338 9.983   1.00 24.93 .
ATOM   1647 O   O   . SER SER A A 58  58  . 6.398   28.535 9.606   1.00 20.82 .
ATOM   1648 CB  CB  . SER SER A A 58  58  . 4.413   28.842 12.270  1.00 22.06 .
ATOM   1649 OG  OG  . SER SER A A 58  58  . 5.361   29.885 12.405  1.00 28.40 .
ATOM   1650 N   N   . VAL VAL A A 59  59  . 4.166   28.628 9.249   1.00 22.74 .
ATOM   1651 CA  CA  . VAL VAL A A 59  59  . 4.243   29.243 7.928   1.00 18.26 .
ATOM   1652 C   C   . VAL VAL A A 59  59  . 3.232   30.376 7.946   1.00 17.40 .
ATOM   1653 O   O   . VAL VAL A A 59  59  . 2.099   30.191 8.382   1.00 22.30 .
ATOM   1654 CB  CB  . VAL VAL A A 59  59  . 3.846   28.258 6.809   1.00 19.67 .
ATOM   1655 CG1 CG1 . VAL VAL A A 59  59  . 3.731   29.000 5.474   1.00 13.08 .
ATOM   1656 CG2 CG2 . VAL VAL A A 59  59  . 4.880   27.144 6.707   1.00 18.01 .
ATOM   1657 N   N   . THR THR A A 60  60  . 3.640   31.550 7.485   1.00 19.62 .
ATOM   1658 CA  CA  . THR THR A A 60  60  . 2.747   32.703 7.460   1.00 17.41 .
ATOM   1659 C   C   . THR THR A A 60  60  . 2.898   33.460 6.147   1.00 16.60 .
ATOM   1660 O   O   . THR THR A A 60  60  . 3.578   33.006 5.224   1.00 17.67 .
ATOM   1661 CB  CB  . THR THR A A 60  60  . 3.094   33.698 8.583   1.00 22.94 .
ATOM   1662 OG1 OG1 . THR THR A A 60  60  . 4.344   34.338 8.270   1.00 15.08 .
ATOM   1663 CG2 CG2 . THR THR A A 60  60  . 3.217   32.973 9.931   1.00 17.23 .
ATOM   1664 N   N   . ALA ALA A A 61  61  . 2.250   34.620 6.076   1.00 16.39 .
ATOM   1665 CA  CA  . ALA ALA A A 61  61  . 2.349   35.490 4.916   1.00 16.69 .
ATOM   1666 C   C   . ALA ALA A A 61  61  . 2.621   36.867 5.505   1.00 19.41 .
ATOM   1667 O   O   . ALA ALA A A 61  61  . 1.900   37.319 6.400   1.00 19.59 .
ATOM   1668 CB  CB  . ALA ALA A A 61  61  . 1.057   35.489 4.120   1.00 20.71 .
ATOM   1669 N   N   . VAL VAL A A 62  62  . 3.653   37.535 5.001   1.00 14.96 .
ATOM   1670 CA  CA  . VAL VAL A A 62  62  . 4.034   38.841 5.523   1.00 21.07 .
ATOM   1671 C   C   . VAL VAL A A 62  62  . 3.862   40.028 4.583   1.00 16.54 .
ATOM   1672 O   O   . VAL VAL A A 62  62  . 4.153   39.946 3.392   1.00 18.70 .
ATOM   1673 CB  CB  . VAL VAL A A 62  62  . 5.505   38.815 5.983   1.00 16.05 .
ATOM   1674 CG1 CG1 . VAL VAL A A 62  62  . 5.661   37.832 7.133   1.00 22.34 .
ATOM   1675 CG2 CG2 . VAL VAL A A 62  62  . 6.406   38.402 4.819   1.00 15.79 .
ATOM   1676 N   N   . LYS LYS A A 63  63  . 3.390   41.135 5.146   1.00 25.55 .
ATOM   1677 CA  CA  . LYS LYS A A 63  63  . 3.211   42.378 4.410   1.00 31.01 .
ATOM   1678 C   C   . LYS LYS A A 63  63  . 4.505   43.142 4.673   1.00 37.98 .
ATOM   1679 O   O   . LYS LYS A A 63  63  . 4.509   44.227 5.247   1.00 50.47 .
ATOM   1680 CB  CB  . LYS LYS A A 63  63  . 1.988   43.133 4.951   1.00 35.64 .
ATOM   1681 CG  CG  . LYS LYS A A 63  63  . 1.866   44.591 4.517   1.00 50.20 .
ATOM   1682 CD  CD  . LYS LYS A A 63  63  . 1.973   44.773 3.010   1.00 55.71 .
ATOM   1683 CE  CE  . LYS LYS A A 63  63  . 1.890   46.253 2.637   1.00 60.22 .
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ATOM   1684 NZ   NZ   . LYS LYS A A 63  63  . 2.930   47.074 3.335   1.00 52.14 .
ATOM   1685 N    N    . ASN ASN A A 64  64  . 5.611   42.530 4.265   1.00 49.82 .
ATOM   1686 CA   CA   . ASN ASN A A 64  64  . 6.945   43.086 4.448   1.00 49.83 .
ATOM   1687 C    C    . ASN ASN A A 64  64  . 7.827   42.662 3.283   1.00 53.43 .
ATOM   1688 O    O    . ASN ASN A A 64  64  . 8.482   41.617 3.328   1.00 52.17 .
ATOM   1689 CB   CB   . ASN ASN A A 64  64  . 7.543   42.582 5.766   1.00 58.88 .
ATOM   1690 CG   CG   . ASN ASN A A 64  64  . 9.058   42.555 5.747   1.00 61.79 .
ATOM   1691 OD1  OD1  . ASN ASN A A 64  64  . 9.705   43.571 5.501   1.00 69.82 .
ATOM   1692 ND2  ND2  . ASN ASN A A 64  64  . 9.632   41.384 6.004   1.00 63.98 .
ATOM   1693 N    N    . SER SER A A 65  65  . 7.839   43.490 2.245   1.00 52.80 .
ATOM   1694 CA   CA   . SER SER A A 65  65  . 8.616   43.230 1.039   1.00 62.28 .
ATOM   1695 C    C    . SER SER A A 65  65  . 10.060  42.795 1.279   1.00 65.05 .
ATOM   1696 O    O    . SER SER A A 65  65  . 10.601  42.942 2.377   1.00 66.34 .
ATOM   1697 CB   CB   . SER SER A A 65  65  . 8.612   44.473 0.145   1.00 59.10 .
ATOM   1698 OG   OG   . SER SER A A 65  65  . 9.398   44.262 -1.013  1.00 64.88 .
ATOM   1699 N    N    . ASP ASP A A 66  66  . 10.665  42.255 0.225   1.00 65.79 .
ATOM   1700 CA   CA   . ASP ASP A A 66  66  . 12.049  41.796 0.232   1.00 68.45 .
ATOM   1701 C    C    . ASP ASP A A 66  66  . 12.596  41.382 1.604   1.00 67.58 .
ATOM   1702 O    O    . ASP ASP A A 66  66  . 13.089  42.217 2.368   1.00 67.39 .
ATOM   1703 CB   CB   . ASP ASP A A 66  66  . 12.945  42.886 -0.376  1.00 72.29 .
ATOM   1704 CG   CG   . ASP ASP A A 66  66  . 13.795  42.374 -1.531  1.00 75.91 .
ATOM   1705 OD1  OD1  . ASP ASP A A 66  66  . 14.539  43.180 -2.134  1.00 78.05 .
ATOM   1706 OD2  OD2  . ASP ASP A A 66  66  . 13.721  41.165 -1.839  1.00 79.93 .
ATOM   1707 N    N    . ASP ASP A A 67  67  . 12.502  40.089 1.905   1.00 58.48 .
ATOM   1708 CA   CA   . ASP ASP A A 67  67  . 13.001  39.529 3.163   1.00 48.34 .
ATOM   1709 C    C    . ASP ASP A A 67  67  . 13.495  38.118 2.842   1.00 41.75 .
ATOM   1710 O    O    . ASP ASP A A 67  67  . 12.782  37.329 2.217   1.00 31.83 .
ATOM   1711 CB   CB   . ASP ASP A A 67  67  . 11.886  39.478 4.219   1.00 57.14 .
ATOM   1712 CG   CG   . ASP ASP A A 67  67  . 12.370  38.947 5.570   1.00 56.40 .
ATOM   1713 OD1  OD1  . ASP ASP A A 67  67  . 11.571  38.941 6.532   1.00 59.39 .
ATOM   1714 OD2  OD2  . ASP ASP A A 67  67  . 13.542  38.533 5.675   1.00 54.73 .
ATOM   1715 N    N    . ILE ILE A A 68  68  . 14.718  37.814 3.262   1.00 32.35 .
ATOM   1716 CA   CA   . ILE ILE A A 68  68  . 15.319  36.514 3.000   1.00 25.91 .
ATOM   1717 C    C    . ILE ILE A A 68  68  . 14.474  35.329 3.481   1.00 19.03 .
ATOM   1718 O    O    . ILE ILE A A 68  68  . 14.504  34.258 2.875   1.00 19.16 .
ATOM   1719 CB   CB   . ILE ILE A A 68  68  . 16.742  36.432 3.622   1.00 34.74 .
ATOM   1720 CG1  CG1  . ILE ILE A A 68  68  . 17.419  35.117 3.216   1.00 33.93 .
ATOM   1721 CG2  CG2  . ILE ILE A A 68  68  . 16.663  36.551 5.136   1.00 32.14 .
ATOM   1722 CD1  CD1  . ILE ILE A A 68  68  . 17.638  34.969 1.717   1.00 39.73 .
ATOM   1723 N    N    . ASN ASN A A 69  69  . 13.725  35.515 4.563   1.00 15.02 .
ATOM   1724 CA   CA   . ASN ASN A A 69  69  . 12.884  34.448 5.080   1.00 19.43 .
ATOM   1725 C    C    . ASN ASN A A 69  69  . 11.626  34.258 4.227   1.00 18.39 .
ATOM   1726 O    O    . ASN ASN A A 69  69  . 10.786  33.424 4.553   1.00 15.94 .
ATOM   1727 CB   CB   . ASN ASN A A 69  69  . 12.480  34.724 6.537   1.00 16.04 .
ATOM   1728 CG   CG   . ASN ASN A A 69  69  . 13.662  34.639 7.498   1.00 22.48 .
ATOM   1729 OD1  OD1  . ASN ASN A A 69  69  . 14.339  33.615 7.576   1.00 24.04 .
ATOM   1730 ND2  ND2  . ASN ASN A A 69  69  . 13.917  35.719 8.221   1.00 25.64 .
ATOM   1731 N    N    . SER SER A A 70  70  . 11.499  35.032 3.148   1.00 10.93 .
ATOM   1732 CA   CA   . SER SER A A 70  70  . 10.341  34.913 2.261   1.00 16.19 .
ATOM   1733 C    C    . SER SER A A 70  70  . 10.726  34.224 0.962   1.00 21.33 .
ATOM   1734 O    O    . SER SER A A 70  70  . 9.923   34.160 0.022   1.00 17.16 .
ATOM   1735 CB   CB   . SER SER A A 70  70  . 9.747   36.289 1.931   1.00 22.74 .
ATOM   1736 OG   OG   . SER SER A A 70  70  . 9.201   36.900 3.089   1.00 20.35 .
ATOM   1737 N    N    . HIS HIS A A 71  71  . 11.960  33.729 0.898   1.00 15.32 .
ATOM   1738 CA   CA   . HIS HIS A A 71  71  . 12.420  33.045 -0.303  1.00 15.60 .
ATOM   1739 C    C    . HIS HIS A A 71  71  . 12.238  31.545 -0.213  1.00 17.97 .
ATOM   1740 O    O    . HIS HIS A A 71  71  . 12.542  30.928 0.821   1.00 15.73 .
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ATOM   1741 CB  CB  . HIS HIS A A 71  71  . 13.884  33.375 -0.582  1.00 20.63 .
ATOM   1742 CG  CG  . HIS HIS A A 71  71  . 14.084  34.773 -1.069  1.00 29.15 .
ATOM   1743 ND1 ND1 . HIS HIS A A 71  71  . 14.055  35.862 -0.228  1.00 34.25 .
ATOM   1744 CD2 CD2 . HIS HIS A A 71  71  . 14.219  35.268 -2.321  1.00 34.44 .
ATOM   1745 CE1 CE1 . HIS HIS A A 71  71  . 14.160  36.969 -0.941  1.00 32.66 .
ATOM   1746 NE2 NE2 . HIS HIS A A 71  71  . 14.259  36.637 -2.215  1.00 36.38 .
ATOM   1747 N   N   . TRP TRP A A 72  72  . 11.731  30.976 -1.306  1.00 12.16 .
ATOM   1748 CA  CA  . TRP TRP A A 72  72  . 11.481  29.546 -1.408  1.00 11.86 .
ATOM   1749 C   C   . TRP TRP A A 72  72  . 12.102  28.986 -2.676  1.00 14.14 .
ATOM   1750 O   O   . TRP TRP A A 72  72  . 12.168  29.664 -3.701  1.00 21.98 .
ATOM   1751 CB  CB  . TRP TRP A A 72  72  . 9.979   29.267 -1.441  1.00 17.14 .
ATOM   1752 CG  CG  . TRP TRP A A 72  72  . 9.243   29.782 -0.266  1.00 10.73 .
ATOM   1753 CD1 CD1 . TRP TRP A A 72  72  . 8.842   31.068 -0.046  1.00 11.99 .
ATOM   1754 CD2 CD2 . TRP TRP A A 72  72  . 8.818   29.025 0.864   1.00 6.42  .
ATOM   1755 NE1 NE1 . TRP TRP A A 72  72  . 8.189   31.157 1.162   1.00 13.45 .
ATOM   1756 CE2 CE2 . TRP TRP A A 72  72  . 8.160   29.915 1.739   1.00 9.57  .
ATOM   1757 CE3 CE3 . TRP TRP A A 72  72  . 8.928   27.671 1.226   1.00 15.25 .
ATOM   1758 CZ2 CZ2 . TRP TRP A A 72  72  . 7.610   29.500 2.954   1.00 12.29 .
ATOM   1759 CZ3 CZ3 . TRP TRP A A 72  72  . 8.382   27.256 2.433   1.00 10.32 .
ATOM   1760 CH2 CH2 . TRP TRP A A 72  72  . 7.729   28.171 3.283   1.00 12.75 .
ATOM   1761 N   N   . GLN GLN A A 73  73  . 12.548  27.739 -2.616  1.00 17.71 .
ATOM   1762 CA  CA  . GLN GLN A A 73  73  . 13.143  27.116 -3.783  1.00 14.84 .
ATOM   1763 C   C   . GLN GLN A A 73  73  . 12.281  25.983 -4.296  1.00 12.80 .
ATOM   1764 O   O   . GLN GLN A A 73  73  . 11.880  25.091 -3.547  1.00 14.15 .
ATOM   1765 CB  CB  . GLN GLN A A 73  73  . 14.540  26.589 -3.457  1.00 17.45 .
ATOM   1766 CG  CG  . GLN GLN A A 73  73  . 15.230  25.952 -4.653  1.00 13.96 .
ATOM   1767 CD  CD  . GLN GLN A A 73  73  . 16.701  25.725 -4.395  1.00 24.58 .
ATOM   1768 OE1 OE1 . GLN GLN A A 73  73  . 17.093  25.381 -3.284  1.00 19.66 .
ATOM   1769 NE2 NE2 . GLN GLN A A 73  73  . 17.523  25.907 -5.424  1.00 26.83 .
ATOM   1770 N   N   . ILE ILE A A 74  74  . 11.992  26.024 -5.586  1.00 15.05 .
ATOM   1771 CA  CA  . ILE ILE A A 74  74  . 11.179  24.997 -6.211  1.00 13.28 .
ATOM   1772 C   C   . ILE ILE A A 74  74  . 11.979  23.720 -6.419  1.00 16.62 .
ATOM   1773 O   O   . ILE ILE A A 74  74  . 13.047  23.740 -7.034  1.00 18.18 .
ATOM   1774 CB  CB  . ILE ILE A A 74  74  . 10.638  25.480 -7.567  1.00 20.34 .
ATOM   1775 CG1 CG1 . ILE ILE A A 74  74  . 9.788   26.736 -7.353  1.00 16.93 .
ATOM   1776 CG2 CG2 . ILE ILE A A 74  74  . 9.809   24.379 -8.220  1.00 25.84 .
ATOM   1777 CD1 CD1 . ILE ILE A A 74  74  . 9.353   27.381 -8.615  1.00 26.87 .
ATOM   1778 N   N   . PHE PHE A A 75  75  . 11.453  22.618 -5.882  1.00 18.79 .
ATOM   1779 CA  CA  . PHE PHE A A 75  75  . 12.066  21.297 -5.997  1.00 17.95 .
ATOM   1780 C   C   . PHE PHE A A 75  75  . 11.132  20.358 -6.755  1.00 22.55 .
ATOM   1781 O   O   . PHE PHE A A 75  75  . 9.917   20.557 -6.760  1.00 16.92 .
ATOM   1782 CB  CB  . PHE PHE A A 75  75  . 12.306  20.688 -4.611  1.00 18.01 .
ATOM   1783 CG  CG  . PHE PHE A A 75  75  . 13.669  20.966 -4.045  1.00 19.86 .
ATOM   1784 CD1 CD1 . PHE PHE A A 75  75  . 14.062  22.259 -3.737  1.00 17.42 .
ATOM   1785 CD2 CD2 . PHE PHE A A 75  75  . 14.559  19.923 -3.814  1.00 28.23 .
ATOM   1786 CE1 CE1 . PHE PHE A A 75  75  . 15.324  22.515 -3.201  1.00 22.82 .
ATOM   1787 CE2 CE2 . PHE PHE A A 75  75  . 15.831  20.168 -3.277  1.00 28.46 .
ATOM   1788 CZ  CZ  . PHE PHE A A 75  75  . 16.208  21.466 -2.971  1.00 23.30 .
ATOM   1789 N   N   . PRO PRO A A 76  76  . 11.685  19.319 -7.405  1.00 25.91 .
ATOM   1790 CA  CA  . PRO PRO A A 76  76  . 10.803  18.394 -8.127  1.00 24.76 .
ATOM   1791 C   C   . PRO PRO A A 76  76  . 10.022  17.632 -7.062  1.00 19.25 .
ATOM   1792 O   O   . PRO PRO A A 76  76  . 10.395  17.667 -5.886  1.00 20.25 .
ATOM   1793 CB  CB  . PRO PRO A A 76  76  . 11.784  17.507 -8.892  1.00 25.61 .
ATOM   1794 CG  CG  . PRO PRO A A 76  76  . 12.986  17.477 -7.983  1.00 25.49 .
ATOM   1795 CD  CD  . PRO PRO A A 76  76  . 13.100  18.921 -7.541  1.00 23.57 .
ATOM   1796 N   N   . ALA ALA A A 77  77  . 8.940   16.966 -7.457  1.00 16.75 .
ATOM   1797 CA  CA  . ALA ALA A A 77  77  . 8.121   16.210 -6.512  1.00 23.14 .
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ATOM   1798 C   C   . ALA ALA A A 77  77  . 8.968   15.168 -5.781  1.00 21.34 .
ATOM   1799 O   O   . ALA ALA A A 77  77  . 10.038  14.793 -6.261  1.00 19.00 .
ATOM   1800 CB  CB  . ALA ALA A A 77  77  . 6.970   15.535 -7.248  1.00 19.46 .
ATOM   1801 N   N   . LEU LEU A A 78  78  . 8.489   14.704 -4.629  1.00 19.96 .
ATOM   1802 CA  CA  . LEU LEU A A 78  78  . 9.232   13.719 -3.836  1.00 28.31 .
ATOM   1803 C   C   . LEU LEU A A 78  78  . 9.770   12.550 -4.648  1.00 32.85 .
ATOM   1804 O   O   . LEU LEU A A 78  78  . 10.936  12.175 -4.509  1.00 35.59 .
ATOM   1805 CB  CB  . LEU LEU A A 78  78  . 8.375   13.165 -2.695  1.00 23.94 .
ATOM   1806 CG  CG  . LEU LEU A A 78  78  . 9.087   12.090 -1.857  1.00 28.15 .
ATOM   1807 CD1 CD1 . LEU LEU A A 78  78  . 10.256  12.715 -1.112  1.00 34.86 .
ATOM   1808 CD2 CD2 . LEU LEU A A 78  78  . 8.116   11.461 -0.871  1.00 37.55 .
ATOM   1809 N   N   . ASN ASN A A 79  79  . 8.934   11.967 -5.497  1.00 29.54 .
ATOM   1810 CA  CA  . ASN ASN A A 79  79  . 9.386   10.825 -6.288  1.00 40.63 .
ATOM   1811 C   C   . ASN ASN A A 79  79  . 9.692   11.159 -7.745  1.00 42.82 .
ATOM   1812 O   O   . ASN ASN A A 79  79  . 9.547   10.319 -8.635  1.00 39.17 .
ATOM   1813 CB  CB  . ASN ASN A A 79  79  . 8.355   9.697  -6.200  1.00 47.65 .
ATOM   1814 CG  CG  . ASN ASN A A 79  79  . 8.190   9.177  -4.778  1.00 54.90 .
ATOM   1815 OD1 OD1 . ASN ASN A A 79  79  . 9.164   8.766  -4.138  1.00 51.31 .
ATOM   1816 ND2 ND2 . ASN ASN A A 79  79  . 6.959   9.197  -4.276  1.00 46.54 .
ATOM   1817 N   N   . ALA ALA A A 80  80  . 10.133  12.391 -7.975  1.00 33.86 .
ATOM   1818 CA  CA  . ALA ALA A A 80  80  . 10.478  12.850 -9.310  1.00 33.05 .
ATOM   1819 C   C   . ALA ALA A A 80  80  . 11.880  13.436 -9.284  1.00 29.84 .
ATOM   1820 O   O   . ALA ALA A A 80  80  . 12.534  13.467 -8.245  1.00 30.27 .
ATOM   1821 CB  CB  . ALA ALA A A 80  80  . 9.483   13.904 -9.777  1.00 23.84 .
ATOM   1822 N   N   . LYS LYS A A 81  81  . 12.329  13.907 -10.439 1.00 32.52 .
ATOM   1823 CA  CA  . LYS LYS A A 81  81  . 13.643  14.513 -10.571 1.00 37.41 .
ATOM   1824 C   C   . LYS LYS A A 81  81  . 13.562  15.551 -11.685 1.00 34.14 .
ATOM   1825 O   O   . LYS LYS A A 81  81  . 12.721  15.450 -12.576 1.00 34.73 .
ATOM   1826 CB  CB  . LYS LYS A A 81  81  . 14.682  13.444 -10.916 1.00 42.02 .
ATOM   1827 CG  CG  . LYS LYS A A 81  81  . 14.334  12.647 -12.162 1.00 54.63 .
ATOM   1828 CD  CD  . LYS LYS A A 81  81  . 15.303  11.493 -12.388 1.00 61.26 .
ATOM   1829 CE  CE  . LYS LYS A A 81  81  . 14.849  10.624 -13.553 1.00 58.14 .
ATOM   1830 NZ  NZ  . LYS LYS A A 81  81  . 15.716  9.425  -13.721 1.00 60.61 .
ATOM   1831 N   N   . CYS CYS A A 82  82  . 14.429  16.552 -11.634 1.00 26.46 .
ATOM   1832 CA  CA  . CYS CYS A A 82  82  . 14.418  17.587 -12.651 1.00 28.66 .
ATOM   1833 C   C   . CYS CYS A A 82  82  . 15.801  18.190 -12.797 1.00 28.44 .
ATOM   1834 O   O   . CYS CYS A A 82  82  . 16.691  17.931 -11.994 1.00 38.48 .
ATOM   1835 CB  CB  . CYS CYS A A 82  82  . 13.449  18.693 -12.255 1.00 35.74 .
ATOM   1836 SG  SG  . CYS CYS A A 82  82  . 14.131  19.728 -10.921 1.00 39.46 .
ATOM   1837 N   N   . ASN ASN A A 83  83  . 15.968  19.011 -13.824 1.00 36.00 .
ATOM   1838 CA  CA  . ASN ASN A A 83  83  . 17.231  19.687 -14.076 1.00 37.06 .
ATOM   1839 C   C   . ASN ASN A A 83  83  . 17.028  21.187 -13.904 1.00 40.07 .
ATOM   1840 O   O   . ASN ASN A A 83  83  . 16.051  21.749 -14.402 1.00 35.61 .
ATOM   1841 CB  CB  . ASN ASN A A 83  83  . 17.704  19.394 -15.498 1.00 44.70 .
ATOM   1842 CG  CG  . ASN ASN A A 83  83  . 18.115  17.950 -15.686 1.00 44.20 .
ATOM   1843 OD1 OD1 . ASN ASN A A 83  83  . 17.957  17.386 -16.765 1.00 48.26 .
ATOM   1844 ND2 ND2 . ASN ASN A A 83  83  . 18.658  17.348 -14.637 1.00 0.00  .
ATOM   1845 N   N   . ARG ARG A A 84  84  . 17.940  21.835 -13.189 1.00 36.76 .
ATOM   1846 CA  CA  . ARG ARG A A 84  84  . 17.839  23.272 -12.986 1.00 36.21 .
ATOM   1847 C   C   . ARG ARG A A 84  84  . 17.563  23.901 -14.344 1.00 40.03 .
ATOM   1848 O   O   . ARG ARG A A 84  84  . 18.228  23.579 -15.329 1.00 42.96 .
ATOM   1849 CB  CB  . ARG ARG A A 84  84  . 19.144  23.825 -12.400 1.00 41.12 .
ATOM   1850 CG  CG  . ARG ARG A A 84  84  . 20.238  22.782 -12.202 1.00 60.84 .
ATOM   1851 CD  CD  . ARG ARG A A 84  84  . 20.753  22.251 -13.541 1.00 69.11 .
ATOM   1852 NE  NE  . ARG ARG A A 84  84  . 21.477  20.988 -13.400 1.00 63.86 .
ATOM   1853 CZ  CZ  . ARG ARG A A 84  84  . 21.979  20.299 -14.420 1.00 65.32 .
ATOM   1854 NH1 NH1 . ARG ARG A A 84  84  . 21.840  20.752 -15.661 1.00 55.45 .
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ATOM   1855 NH2 NH2 . ARG ARG A A 84  84  . 22.611  19.152 -14.203 1.00 64.96 .
ATOM   1856 N   N   . GLY GLY A A 85  85  . 16.573  24.784 -14.405 1.00 31.47 .
ATOM   1857 CA  CA  . GLY GLY A A 85  85  . 16.245  25.415 -15.668 1.00 35.95 .
ATOM   1858 C   C   . GLY GLY A A 85  85  . 15.067  24.751 -16.359 1.00 38.44 .
ATOM   1859 O   O   . GLY GLY A A 85  85  . 14.492  25.311 -17.295 1.00 40.28 .
ATOM   1860 N   N   . ASP ASP A A 86  86  . 14.704  23.554 -15.908 1.00 32.62 .
ATOM   1861 CA  CA  . ASP ASP A A 86  86  . 13.573  22.841 -16.494 1.00 33.72 .
ATOM   1862 C   C   . ASP ASP A A 86  86  . 12.277  23.584 -16.207 1.00 34.44 .
ATOM   1863 O   O   . ASP ASP A A 86  86  . 12.047  24.039 -15.091 1.00 32.61 .
ATOM   1864 CB  CB  . ASP ASP A A 86  86  . 13.465  21.419 -15.931 1.00 31.68 .
ATOM   1865 CG  CG  . ASP ASP A A 86  86  . 14.491  20.478 -16.523 1.00 38.31 .
ATOM   1866 OD1 OD1 . ASP ASP A A 86  86  . 15.213  20.901 -17.451 1.00 28.40 .
ATOM   1867 OD2 OD2 . ASP ASP A A 86  86  . 14.568  19.316 -16.067 1.00 30.96 .
ATOM   1868 N   N   . ALA ALA A A 87  87  . 11.435  23.709 -17.224 1.00 32.83 .
ATOM   1869 CA  CA  . ALA ALA A A 87  87  . 10.160  24.378 -17.060 1.00 33.11 .
ATOM   1870 C   C   . ALA ALA A A 87  87  . 9.266   23.465 -16.245 1.00 27.79 .
ATOM   1871 O   O   . ALA ALA A A 87  87  . 9.398   22.244 -16.299 1.00 25.57 .
ATOM   1872 CB  CB  . ALA ALA A A 87  87  . 9.530   24.648 -18.420 1.00 34.78 .
ATOM   1873 N   N   . ILE ILE A A 88  88  . 8.371   24.055 -15.465 1.00 33.57 .
ATOM   1874 CA  CA  . ILE ILE A A 88  88  . 7.445   23.262 -14.674 1.00 29.69 .
ATOM   1875 C   C   . ILE ILE A A 88  88  . 6.117   23.309 -15.408 1.00 34.02 .
ATOM   1876 O   O   . ILE ILE A A 88  88  . 5.616   24.388 -15.733 1.00 27.55 .
ATOM   1877 CB  CB  . ILE ILE A A 88  88  . 7.271   23.831 -13.260 1.00 30.67 .
ATOM   1878 CG1 CG1 . ILE ILE A A 88  88  . 8.607   23.756 -12.514 1.00 34.91 .
ATOM   1879 CG2 CG2 . ILE ILE A A 88  88  . 6.190   23.047 -12.510 1.00 23.27 .
ATOM   1880 CD1 CD1 . ILE ILE A A 88  88  . 8.595   24.460 -11.196 1.00 38.11 .
ATOM   1881 N   N   . LYS LYS A A 89  89  . 5.558   22.137 -15.684 1.00 28.62 .
ATOM   1882 CA  CA  . LYS LYS A A 89  89  . 4.295   22.063 -16.395 1.00 32.16 .
ATOM   1883 C   C   . LYS LYS A A 89  89  . 3.146   22.147 -15.403 1.00 29.68 .
ATOM   1884 O   O   . LYS LYS A A 89  89  . 3.270   21.673 -14.274 1.00 23.85 .
ATOM   1885 CB  CB  . LYS LYS A A 89  89  . 4.216   20.748 -17.183 1.00 37.12 .
ATOM   1886 CG  CG  . LYS LYS A A 89  89  . 5.308   20.570 -18.252 1.00 37.96 .
ATOM   1887 CD  CD  . LYS LYS A A 89  89  . 6.674   20.298 -17.625 1.00 45.80 .
ATOM   1888 CE  CE  . LYS LYS A A 89  89  . 7.767   20.151 -18.678 1.00 55.66 .
ATOM   1889 NZ  NZ  . LYS LYS A A 89  89  . 9.095   19.841 -18.070 1.00 46.12 .
ATOM   1890 N   N   . CYS CYS A A 90  90  . 2.033   22.754 -15.816 1.00 26.67 .
ATOM   1891 CA  CA  . CYS CYS A A 90  90  . 0.869   22.867 -14.939 1.00 23.14 .
ATOM   1892 C   C   . CYS CYS A A 90  90  . 0.326   21.470 -14.695 1.00 22.14 .
ATOM   1893 O   O   . CYS CYS A A 90  90  . 0.307   20.637 -15.599 1.00 26.09 .
ATOM   1894 CB  CB  . CYS CYS A A 90  90  . -0.208  23.775 -15.565 1.00 20.81 .
ATOM   1895 SG  SG  . CYS CYS A A 90  90  . 0.459   25.423 -15.967 1.00 32.21 .
ATOM   1896 N   N   . GLY GLY A A 91  91  . -0.101  21.212 -13.466 1.00 19.38 .
ATOM   1897 CA  CA  . GLY GLY A A 91  91  . -0.617  19.906 -13.128 1.00 18.94 .
ATOM   1898 C   C   . GLY GLY A A 91  91  . 0.451   19.055 -12.462 1.00 21.25 .
ATOM   1899 O   O   . GLY GLY A A 91  91  . 0.139   18.089 -11.768 1.00 24.01 .
ATOM   1900 N   N   . ASP ASP A A 92  92  . 1.716   19.397 -12.679 1.00 23.40 .
ATOM   1901 CA  CA  . ASP ASP A A 92  92  . 2.805   18.647 -12.057 1.00 27.21 .
ATOM   1902 C   C   . ASP ASP A A 92  92  . 2.929   19.006 -10.581 1.00 24.19 .
ATOM   1903 O   O   . ASP ASP A A 92  92  . 2.538   20.096 -10.164 1.00 19.26 .
ATOM   1904 CB  CB  . ASP ASP A A 92  92  . 4.135   18.940 -12.763 1.00 29.22 .
ATOM   1905 CG  CG  . ASP ASP A A 92  92  . 4.383   18.023 -13.950 1.00 37.68 .
ATOM   1906 OD1 OD1 . ASP ASP A A 92  92  . 5.420   18.201 -14.621 1.00 46.45 .
ATOM   1907 OD2 OD2 . ASP ASP A A 92  92  . 3.550   17.125 -14.207 1.00 36.00 .
ATOM   1908 N   N   . LYS LYS A A 93  93  . 3.458   18.077 -9.792  1.00 19.99 .
ATOM   1909 CA  CA  . LYS LYS A A 93  93  . 3.656   18.309 -8.371  1.00 20.83 .
ATOM   1910 C   C   . LYS LYS A A 93  93  . 5.057   18.863 -8.147  1.00 23.59 .
ATOM   1911 O   O   . LYS LYS A A 93  93  . 5.973   18.596 -8.929  1.00 16.09 .
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ATOM   1912 CB  CB  . LYS LYS A A 93  93  . 3.483   17.006 -7.578  1.00 21.63 .
ATOM   1913 CG  CG  . LYS LYS A A 93  93  . 2.034   16.549 -7.484  1.00 30.33 .
ATOM   1914 CD  CD  . LYS LYS A A 93  93  . 1.861   15.336 -6.596  1.00 34.71 .
ATOM   1915 CE  CE  . LYS LYS A A 93  93  . 2.412   14.082 -7.244  1.00 38.03 .
ATOM   1916 NZ  NZ  . LYS LYS A A 93  93  . 2.061   12.880 -6.440  1.00 44.92 .
ATOM   1917 N   N   . ILE ILE A A 94  94  . 5.205   19.670 -7.103  1.00 18.64 .
ATOM   1918 CA  CA  . ILE ILE A A 94  94  . 6.497   20.235 -6.744  1.00 17.58 .
ATOM   1919 C   C   . ILE ILE A A 94  94  . 6.592   20.227 -5.234  1.00 14.82 .
ATOM   1920 O   O   . ILE ILE A A 94  94  . 5.646   19.861 -4.545  1.00 16.80 .
ATOM   1921 CB  CB  . ILE ILE A A 94  94  . 6.664   21.704 -7.191  1.00 15.70 .
ATOM   1922 CG1 CG1 . ILE ILE A A 94  94  . 5.602   22.588 -6.519  1.00 19.95 .
ATOM   1923 CG2 CG2 . ILE ILE A A 94  94  . 6.596   21.797 -8.696  1.00 21.39 .
ATOM   1924 CD1 CD1 . ILE ILE A A 94  94  . 5.786   24.090 -6.767  1.00 22.79 .
ATOM   1925 N   N   . ARG ARG A A 95  95  . 7.759   20.611 -4.739  1.00 17.95 .
ATOM   1926 CA  CA  . ARG ARG A A 95  95  . 8.006   20.746 -3.320  1.00 15.08 .
ATOM   1927 C   C   . ARG ARG A A 95  95  . 8.598   22.146 -3.202  1.00 13.78 .
ATOM   1928 O   O   . ARG ARG A A 95  95  . 9.250   22.626 -4.131  1.00 17.71 .
ATOM   1929 CB  CB  . ARG ARG A A 95  95  . 9.002   19.690 -2.831  1.00 16.03 .
ATOM   1930 CG  CG  . ARG ARG A A 95  95  . 8.366   18.332 -2.511  1.00 17.90 .
ATOM   1931 CD  CD  . ARG ARG A A 95  95  . 9.407   17.353 -1.983  1.00 18.32 .
ATOM   1932 NE  NE  . ARG ARG A A 95  95  . 10.409  17.079 -3.004  1.00 18.88 .
ATOM   1933 CZ  CZ  . ARG ARG A A 95  95  . 11.615  16.572 -2.769  1.00 23.43 .
ATOM   1934 NH1 NH1 . ARG ARG A A 95  95  . 11.985  16.274 -1.532  1.00 19.98 .
ATOM   1935 NH2 NH2 . ARG ARG A A 95  95  . 12.459  16.381 -3.778  1.00 14.71 .
ATOM   1936 N   N   . LEU LEU A A 96  96  . 8.355   22.815 -2.081  1.00 15.59 .
ATOM   1937 CA  CA  . LEU LEU A A 96  96  . 8.876   24.158 -1.891  1.00 15.09 .
ATOM   1938 C   C   . LEU LEU A A 96  96  . 9.756   24.196 -0.650  1.00 12.27 .
ATOM   1939 O   O   . LEU LEU A A 96  96  . 9.286   23.989 0.469   1.00 13.79 .
ATOM   1940 CB  CB  . LEU LEU A A 96  96  . 7.727   25.167 -1.753  1.00 20.55 .
ATOM   1941 CG  CG  . LEU LEU A A 96  96  . 6.961   25.509 -3.037  1.00 25.97 .
ATOM   1942 CD1 CD1 . LEU LEU A A 96  96  . 5.777   26.423 -2.716  1.00 23.11 .
ATOM   1943 CD2 CD2 . LEU LEU A A 96  96  . 7.908   26.199 -4.025  1.00 18.62 .
ATOM   1944 N   N   . LYS LYS A A 97  97  . 11.038  24.461 -0.867  1.00 15.43 .
ATOM   1945 CA  CA  . LYS LYS A A 97  97  . 12.008  24.523 0.218   1.00 13.62 .
ATOM   1946 C   C   . LYS LYS A A 97  97  . 12.199  25.938 0.767   1.00 13.33 .
ATOM   1947 O   O   . LYS LYS A A 97  97  . 12.454  26.876 0.008   1.00 16.33 .
ATOM   1948 CB  CB  . LYS LYS A A 97  97  . 13.350  23.989 -0.282  1.00 13.69 .
ATOM   1949 CG  CG  . LYS LYS A A 97  97  . 14.394  23.817 0.816   1.00 20.49 .
ATOM   1950 CD  CD  . LYS LYS A A 97  97  . 15.694  23.311 0.231   1.00 23.55 .
ATOM   1951 CE  CE  . LYS LYS A A 97  97  . 16.711  23.006 1.311   1.00 33.61 .
ATOM   1952 NZ  NZ  . LYS LYS A A 97  97  . 17.974  22.541 0.698   1.00 29.81 .
ATOM   1953 N   N   . HIS HIS A A 98  98  . 12.051  26.097 2.076   1.00 10.64 .
ATOM   1954 CA  CA  . HIS HIS A A 98  98  . 12.258  27.396 2.698   1.00 15.52 .
ATOM   1955 C   C   . HIS HIS A A 98  98  . 13.775  27.515 2.835   1.00 21.16 .
ATOM   1956 O   O   . HIS HIS A A 98  98  . 14.398  26.773 3.601   1.00 18.88 .
ATOM   1957 CB  CB  . HIS HIS A A 98  98  . 11.575  27.451 4.060   1.00 14.29 .
ATOM   1958 CG  CG  . HIS HIS A A 98  98  . 11.700  28.777 4.735   1.00 22.33 .
ATOM   1959 ND1 ND1 . HIS HIS A A 98  98  . 12.337  28.937 5.947   1.00 17.69 .
ATOM   1960 CD2 CD2 . HIS HIS A A 98  98  . 11.266  30.009 4.369   1.00 15.86 .
ATOM   1961 CE1 CE1 . HIS HIS A A 98  98  . 12.285  30.211 6.302   1.00 24.14 .
ATOM   1962 NE2 NE2 . HIS HIS A A 98  98  . 11.642  30.882 5.361   1.00 18.82 .
ATOM   1963 N   N   . LEU LEU A A 99  99  . 14.361  28.440 2.082   1.00 18.40 .
ATOM   1964 CA  CA  . LEU LEU A A 99  99  . 15.809  28.606 2.052   1.00 22.07 .
ATOM   1965 C   C   . LEU LEU A A 99  99  . 16.555  28.784 3.359   1.00 27.89 .
ATOM   1966 O   O   . LEU LEU A A 99  99  . 17.542  28.093 3.597   1.00 28.20 .
ATOM   1967 CB  CB  . LEU LEU A A 99  99  . 16.194  29.750 1.110   1.00 27.86 .
ATOM   1968 CG  CG  . LEU LEU A A 99  99  . 16.225  29.443 -0.389  1.00 37.42 .
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ATOM   1969 CD1 CD1 . LEU LEU A A 99  99  . 16.891  30.601 -1.115  1.00 39.90 .
ATOM   1970 CD2 CD2 . LEU LEU A A 99  99  . 16.996  28.153 -0.647  1.00 39.14 .
ATOM   1971 N   N   . THR THR A A 100 100 . 16.093  29.687 4.212   1.00 24.68 .
ATOM   1972 CA  CA  . THR THR A A 100 100 . 16.798  29.937 5.461   1.00 33.21 .
ATOM   1973 C   C   . THR THR A A 100 100 . 16.722  28.850 6.528   1.00 36.80 .
ATOM   1974 O   O   . THR THR A A 100 100 . 17.573  28.815 7.416   1.00 36.81 .
ATOM   1975 CB  CB  . THR THR A A 100 100 . 16.356  31.272 6.087   1.00 27.30 .
ATOM   1976 OG1 OG1 . THR THR A A 100 100 . 14.966  31.212 6.413   1.00 22.45 .
ATOM   1977 CG2 CG2 . THR THR A A 100 100 . 16.590  32.416 5.103   1.00 17.20 .
ATOM   1978 N   N   . THR THR A A 101 101 . 15.727  27.966 6.463   1.00 26.74 .
ATOM   1979 CA  CA  . THR THR A A 101 101 . 15.636  26.910 7.472   1.00 24.64 .
ATOM   1980 C   C   . THR THR A A 101 101 . 15.772  25.507 6.895   1.00 26.22 .
ATOM   1981 O   O   . THR THR A A 101 101 . 15.820  24.530 7.638   1.00 27.05 .
ATOM   1982 CB  CB  . THR THR A A 101 101 . 14.323  26.990 8.266   1.00 23.10 .
ATOM   1983 OG1 OG1 . THR THR A A 101 101 . 13.211  26.742 7.393   1.00 17.86 .
ATOM   1984 CG2 CG2 . THR THR A A 101 101 . 14.182  28.365 8.907   1.00 25.16 .
ATOM   1985 N   N   . GLY GLY A A 102 102 . 15.829  25.414 5.571   1.00 24.79 .
ATOM   1986 CA  CA  . GLY GLY A A 102 102 . 15.984  24.126 4.920   1.00 32.71 .
ATOM   1987 C   C   . GLY GLY A A 102 102 . 14.791  23.188 4.978   1.00 30.03 .
ATOM   1988 O   O   . GLY GLY A A 102 102 . 14.891  22.032 4.561   1.00 2.31  .
ATOM   1989 N   N   . THR THR A A 103 103 . 13.672  23.684 5.494   1.00 17.98 .
ATOM   1990 CA  CA  . THR THR A A 103 103 . 12.443  22.912 5.613   1.00 19.89 .
ATOM   1991 C   C   . THR THR A A 103 103 . 11.633  22.945 4.304   1.00 19.98 .
ATOM   1992 O   O   . THR THR A A 103 103 . 12.005  23.638 3.352   1.00 15.58 .
ATOM   1993 CB  CB  . THR THR A A 103 103 . 11.564  23.487 6.727   1.00 18.62 .
ATOM   1994 OG1 OG1 . THR THR A A 103 103 . 11.433  24.896 6.525   1.00 23.14 .
ATOM   1995 CG2 CG2 . THR THR A A 103 103 . 12.175  23.230 8.103   1.00 22.72 .
ATOM   1996 N   N   . PHE PHE A A 104 104 . 10.527  22.202 4.269   1.00 17.92 .
ATOM   1997 CA  CA  . PHE PHE A A 104 104 . 9.660   22.162 3.091   1.00 18.35 .
ATOM   1998 C   C   . PHE PHE A A 104 104 . 8.216   22.527 3.435   1.00 18.27 .
ATOM   1999 O   O   . PHE PHE A A 104 104 . 7.729   22.209 4.521   1.00 23.05 .
ATOM   2000 CB  CB  . PHE PHE A A 104 104 . 9.677   20.766 2.441   1.00 22.41 .
ATOM   2001 CG  CG  . PHE PHE A A 104 104 . 10.881  20.508 1.571   1.00 21.72 .
ATOM   2002 CD1 CD1 . PHE PHE A A 104 104 . 12.033  19.938 2.100   1.00 18.02 .
ATOM   2003 CD2 CD2 . PHE PHE A A 104 104 . 10.862  20.853 0.220   1.00 15.54 .
ATOM   2004 CE1 CE1 . PHE PHE A A 104 104 . 13.153  19.714 1.301   1.00 15.30 .
ATOM   2005 CE2 CE2 . PHE PHE A A 104 104 . 11.972  20.635 -0.586  1.00 13.76 .
ATOM   2006 CZ  CZ  . PHE PHE A A 104 104 . 13.125  20.061 -0.043  1.00 20.90 .
ATOM   2007 N   N   . LEU LEU A A 105 105 . 7.531   23.203 2.514   1.00 19.81 .
ATOM   2008 CA  CA  . LEU LEU A A 105 105 . 6.138   23.563 2.738   1.00 12.06 .
ATOM   2009 C   C   . LEU LEU A A 105 105 . 5.438   22.218 2.927   1.00 17.67 .
ATOM   2010 O   O   . LEU LEU A A 105 105 . 5.489   21.349 2.049   1.00 19.96 .
ATOM   2011 CB  CB  . LEU LEU A A 105 105 . 5.577   24.275 1.516   1.00 16.62 .
ATOM   2012 CG  CG  . LEU LEU A A 105 105 . 4.212   24.914 1.717   1.00 15.69 .
ATOM   2013 CD1 CD1 . LEU LEU A A 105 105 . 4.312   26.011 2.790   1.00 15.71 .
ATOM   2014 CD2 CD2 . LEU LEU A A 105 105 . 3.731   25.482 0.393   1.00 13.78 .
ATOM   2015 N   N   . HIS HIS A A 106 106 . 4.760   22.064 4.053   1.00 17.70 .
ATOM   2016 CA  CA  . HIS HIS A A 106 106 . 4.142   20.790 4.398   1.00 24.95 .
ATOM   2017 C   C   . HIS HIS A A 106 106 . 2.703   20.897 4.873   1.00 22.73 .
ATOM   2018 O   O   . HIS HIS A A 106 106 . 2.281   21.926 5.403   1.00 22.46 .
ATOM   2019 CB  CB  . HIS HIS A A 106 106 . 4.980   20.163 5.517   1.00 20.61 .
ATOM   2020 CG  CG  . HIS HIS A A 106 106 . 4.663   18.731 5.808   1.00 20.55 .
ATOM   2021 ND1 ND1 . HIS HIS A A 106 106 . 5.308   17.686 5.179   1.00 22.75 .
ATOM   2022 CD2 CD2 . HIS HIS A A 106 106 . 3.828   18.168 6.714   1.00 29.16 .
ATOM   2023 CE1 CE1 . HIS HIS A A 106 106 . 4.889   16.542 5.692   1.00 23.47 .
ATOM   2024 NE2 NE2 . HIS HIS A A 106 106 . 3.992   16.806 6.625   1.00 22.42 .
ATOM   2025 N   N   . SER SER A A 107 107 . 1.942   19.829 4.660   1.00 23.14 .
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ATOM   2026 CA  CA  . SER SER A A 107 107 . 0.577   19.790 5.151   1.00 21.91 .
ATOM   2027 C   C   . SER SER A A 107 107 . 0.251   18.388 5.633   1.00 27.58 .
ATOM   2028 O   O   . SER SER A A 107 107 . 0.973   17.428 5.347   1.00 20.45 .
ATOM   2029 CB  CB  . SER SER A A 107 107 . -0.431  20.235 4.105   1.00 21.58 .
ATOM   2030 OG  OG  . SER SER A A 107 107 . -1.569  20.755 4.782   1.00 20.62 .
ATOM   2031 N   N   . HIS HIS A A 108 108 . -0.845  18.276 6.366   1.00 31.87 .
ATOM   2032 CA  CA  . HIS HIS A A 108 108 . -1.229  17.006 6.952   1.00 33.03 .
ATOM   2033 C   C   . HIS HIS A A 108 108 . -2.559  17.152 7.669   1.00 33.58 .
ATOM   2034 O   O   . HIS HIS A A 108 108 . -3.193  18.205 7.603   1.00 29.61 .
ATOM   2035 CB  CB  . HIS HIS A A 108 108 . -0.138  16.579 7.932   1.00 30.72 .
ATOM   2036 CG  CG  . HIS HIS A A 108 108 . 0.470   17.723 8.686   1.00 28.32 .
ATOM   2037 ND1 ND1 . HIS HIS A A 108 108 . 1.662   17.615 9.368   1.00 27.84 .
ATOM   2038 CD2 CD2 . HIS HIS A A 108 108 . 0.050   18.999 8.864   1.00 29.18 .
ATOM   2039 CE1 CE1 . HIS HIS A A 108 108 . 1.950   18.775 9.932   1.00 32.72 .
ATOM   2040 NE2 NE2 . HIS HIS A A 108 108 . 0.988   19.632 9.641   1.00 22.60 .
ATOM   2041 N   N   . HIS HIS A A 109 109 . -2.973  16.094 8.360   1.00 39.24 .
ATOM   2042 CA  CA  . HIS HIS A A 109 109 . -4.242  16.103 9.068   1.00 43.99 .
ATOM   2043 C   C   . HIS HIS A A 109 109 . -4.240  16.710 10.467  1.00 41.96 .
ATOM   2044 O   O   . HIS HIS A A 109 109 . -4.641  16.063 11.431  1.00 44.14 .
ATOM   2045 CB  CB  . HIS HIS A A 109 109 . -4.819  14.686 9.134   1.00 49.87 .
ATOM   2046 CG  CG  . HIS HIS A A 109 109 . -5.489  14.250 7.869   1.00 61.18 .
ATOM   2047 ND1 ND1 . HIS HIS A A 109 109 . -6.470  14.997 7.253   1.00 65.80 .
ATOM   2048 CD2 CD2 . HIS HIS A A 109 109 . -5.337  13.134 7.116   1.00 68.17 .
ATOM   2049 CE1 CE1 . HIS HIS A A 109 109 . -6.894  14.361 6.176   1.00 66.87 .
ATOM   2050 NE2 NE2 . HIS HIS A A 109 109 . -6.224  13.228 6.070   1.00 68.33 .
ATOM   2051 N   N   . PHE PHE A A 110 110 . -3.777  17.947 10.577  1.00 36.73 .
ATOM   2052 CA  CA  . PHE PHE A A 110 110 . -3.791  18.642 11.854  1.00 38.47 .
ATOM   2053 C   C   . PHE PHE A A 110 110 . -4.641  19.884 11.631  1.00 34.69 .
ATOM   2054 O   O   . PHE PHE A A 110 110 . -4.807  20.331 10.498  1.00 39.03 .
ATOM   2055 CB  CB  . PHE PHE A A 110 110 . -2.376  19.033 12.305  1.00 37.20 .
ATOM   2056 CG  CG  . PHE PHE A A 110 110 . -1.517  17.863 12.708  1.00 36.66 .
ATOM   2057 CD1 CD1 . PHE PHE A A 110 110 . -0.729  17.204 11.773  1.00 38.68 .
ATOM   2058 CD2 CD2 . PHE PHE A A 110 110 . -1.504  17.416 14.026  1.00 40.00 .
ATOM   2059 CE1 CE1 . PHE PHE A A 110 110 . 0.065   16.111 12.144  1.00 49.27 .
ATOM   2060 CE2 CE2 . PHE PHE A A 110 110 . -0.718  16.328 14.409  1.00 42.84 .
ATOM   2061 CZ  CZ  . PHE PHE A A 110 110 . 0.070   15.673 13.465  1.00 47.93 .
ATOM   2062 N   N   . THR THR A A 111 111 . -5.187  20.439 12.704  1.00 40.91 .
ATOM   2063 CA  CA  . THR THR A A 111 111 . -6.035  21.617 12.586  1.00 38.98 .
ATOM   2064 C   C   . THR THR A A 111 111 . -5.197  22.895 12.590  1.00 38.77 .
ATOM   2065 O   O   . THR THR A A 111 111 . -4.199  22.988 13.308  1.00 40.14 .
ATOM   2066 CB  CB  . THR THR A A 111 111 . -7.059  21.661 13.739  1.00 41.08 .
ATOM   2067 OG1 OG1 . THR THR A A 111 111 . -8.103  22.589 13.421  1.00 52.27 .
ATOM   2068 CG2 CG2 . THR THR A A 111 111 . -6.384  22.090 15.033  1.00 43.62 .
ATOM   2069 N   N   . ALA ALA A A 112 112 . -5.600  23.870 11.777  1.00 31.39 .
ATOM   2070 CA  CA  . ALA ALA A A 112 112 . -4.892  25.147 11.689  1.00 37.46 .
ATOM   2071 C   C   . ALA ALA A A 112 112 . -4.776  25.762 13.081  1.00 40.25 .
ATOM   2072 O   O   . ALA ALA A A 112 112 . -5.578  25.461 13.967  1.00 42.76 .
ATOM   2073 CB  CB  . ALA ALA A A 112 112 . -5.637  26.100 10.753  1.00 30.39 .
ATOM   2074 N   N   . PRO PRO A A 113 113 . -3.775  26.634 13.292  1.00 38.42 .
ATOM   2075 CA  CA  . PRO PRO A A 113 113 . -3.566  27.283 14.592  1.00 35.67 .
ATOM   2076 C   C   . PRO PRO A A 113 113 . -4.660  28.247 15.060  1.00 37.03 .
ATOM   2077 O   O   . PRO PRO A A 113 113 . -5.039  28.223 16.226  1.00 40.77 .
ATOM   2078 CB  CB  . PRO PRO A A 113 113 . -2.219  27.976 14.416  1.00 29.83 .
ATOM   2079 CG  CG  . PRO PRO A A 113 113 . -2.207  28.311 12.961  1.00 29.44 .
ATOM   2080 CD  CD  . PRO PRO A A 113 113 . -2.730  27.038 12.336  1.00 36.05 .
ATOM   2081 N   N   . LEU LEU A A 114 114 . -5.161  29.094 14.166  1.00 35.63 .
ATOM   2082 CA  CA  . LEU LEU A A 114 114 . -6.199  30.054 14.543  1.00 36.64 .
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ATOM   2083 C   C   . LEU LEU A A 114 114 . -7.538  29.759 13.885  1.00 42.63 .
ATOM   2084 O   O   . LEU LEU A A 114 114 . -8.577  30.242 14.335  1.00 47.02 .
ATOM   2085 CB  CB  . LEU LEU A A 114 114 . -5.770  31.482 14.186  1.00 38.74 .
ATOM   2086 CG  CG  . LEU LEU A A 114 114 . -4.469  32.010 14.799  1.00 40.77 .
ATOM   2087 CD1 CD1 . LEU LEU A A 114 114 . -4.199  33.410 14.273  1.00 39.69 .
ATOM   2088 CD2 CD2 . LEU LEU A A 114 114 . -4.567  32.014 16.317  1.00 40.79 .
ATOM   2089 N   N   . SER SER A A 115 115 . -7.516  28.975 12.814  1.00 36.08 .
ATOM   2090 CA  CA  . SER SER A A 115 115 . -8.748  28.622 12.118  1.00 41.78 .
ATOM   2091 C   C   . SER SER A A 115 115 . -9.025  27.134 12.273  1.00 42.25 .
ATOM   2092 O   O   . SER SER A A 115 115 . -8.560  26.311 11.485  1.00 39.09 .
ATOM   2093 CB  CB  . SER SER A A 115 115 . -8.648  28.993 10.640  1.00 39.54 .
ATOM   2094 OG  OG  . SER SER A A 115 115 . -8.555  30.397 10.489  1.00 35.31 .
ATOM   2095 N   N   . LYS LYS A A 116 116 . -9.793  26.805 13.306  1.00 45.24 .
ATOM   2096 CA  CA  . LYS LYS A A 116 116 . -10.150 25.427 13.623  1.00 45.06 .
ATOM   2097 C   C   . LYS LYS A A 116 116 . -10.745 24.637 12.459  1.00 42.14 .
ATOM   2098 O   O   . LYS LYS A A 116 116 . -10.520 23.428 12.343  1.00 43.54 .
ATOM   2099 CB  CB  . LYS LYS A A 116 116 . -11.137 25.416 14.792  1.00 48.52 .
ATOM   2100 CG  CG  . LYS LYS A A 116 116 . -10.577 25.958 16.099  1.00 52.57 .
ATOM   2101 CD  CD  . LYS LYS A A 116 116 . -9.574  24.998 16.722  1.00 57.98 .
ATOM   2102 CE  CE  . LYS LYS A A 116 116 . -9.225  25.410 18.150  1.00 61.18 .
ATOM   2103 NZ  NZ  . LYS LYS A A 116 116 . -8.426  24.369 18.872  1.00 58.28 .
ATOM   2104 N   N   . GLN GLN A A 117 117 . -11.506 25.313 11.603  1.00 36.04 .
ATOM   2105 CA  CA  . GLN GLN A A 117 117 . -12.144 24.651 10.471  1.00 38.17 .
ATOM   2106 C   C   . GLN GLN A A 117 117 . -11.180 24.309 9.346   1.00 40.86 .
ATOM   2107 O   O   . GLN GLN A A 117 117 . -11.542 23.597 8.403   1.00 39.81 .
ATOM   2108 CB  CB  . GLN GLN A A 117 117 . -13.268 25.517 9.898   1.00 38.01 .
ATOM   2109 CG  CG  . GLN GLN A A 117 117 . -13.562 26.811 10.649  1.00 49.40 .
ATOM   2110 CD  CD  . GLN GLN A A 117 117 . -12.477 27.862 10.496  1.00 36.04 .
ATOM   2111 OE1 OE1 . GLN GLN A A 117 117 . -11.435 27.796 11.147  1.00 47.61 .
ATOM   2112 NE2 NE2 . GLN GLN A A 117 117 . -12.717 28.838 9.626   1.00 30.21 .
ATOM   2113 N   N   . HIS HIS A A 118 118 . -9.956  24.816 9.436   1.00 35.91 .
ATOM   2114 CA  CA  . HIS HIS A A 118 118 . -8.986  24.558 8.392   1.00 31.16 .
ATOM   2115 C   C   . HIS HIS A A 118 118 . -7.818  23.680 8.799   1.00 35.68 .
ATOM   2116 O   O   . HIS HIS A A 118 118 . -7.559  23.449 9.983   1.00 37.44 .
ATOM   2117 CB  CB  . HIS HIS A A 118 118 . -8.483  25.879 7.805   1.00 34.15 .
ATOM   2118 CG  CG  . HIS HIS A A 118 118 . -9.538  26.639 7.062   1.00 30.13 .
ATOM   2119 ND1 ND1 . HIS HIS A A 118 118 . -10.397 26.033 6.170   1.00 34.64 .
ATOM   2120 CD2 CD2 . HIS HIS A A 118 118 . -9.865  27.952 7.065   1.00 30.49 .
ATOM   2121 CE1 CE1 . HIS HIS A A 118 118 . -11.207 26.939 5.653   1.00 34.06 .
ATOM   2122 NE2 NE2 . HIS HIS A A 118 118 . -10.905 28.113 6.180   1.00 36.14 .
ATOM   2123 N   N   . GLN GLN A A 119 119 . -7.123  23.195 7.779   1.00 32.92 .
ATOM   2124 CA  CA  . GLN GLN A A 119 119 . -5.981  22.314 7.928   1.00 29.72 .
ATOM   2125 C   C   . GLN GLN A A 119 119 . -4.700  23.125 8.113   1.00 32.33 .
ATOM   2126 O   O   . GLN GLN A A 119 119 . -4.536  24.193 7.520   1.00 26.83 .
ATOM   2127 CB  CB  . GLN GLN A A 119 119 . -5.902  21.440 6.681   1.00 33.59 .
ATOM   2128 CG  CG  . GLN GLN A A 119 119 . -4.839  20.371 6.665   1.00 35.49 .
ATOM   2129 CD  CD  . GLN GLN A A 119 119 . -4.978  19.492 5.440   1.00 44.06 .
ATOM   2130 OE1 OE1 . GLN GLN A A 119 119 . -5.926  18.710 5.330   1.00 45.61 .
ATOM   2131 NE2 NE2 . GLN GLN A A 119 119 . -4.051  19.631 4.499   1.00 27.33 .
ATOM   2132 N   N   . GLU GLU A A 120 120 . -3.801  22.607 8.943   1.00 27.21 .
ATOM   2133 CA  CA  . GLU GLU A A 120 120 . -2.532  23.262 9.239   1.00 31.98 .
ATOM   2134 C   C   . GLU GLU A A 120 120 . -1.547  23.161 8.081   1.00 33.12 .
ATOM   2135 O   O   . GLU GLU A A 120 120 . -1.504  22.158 7.367   1.00 27.43 .
ATOM   2136 CB  CB  . GLU GLU A A 120 120 . -1.888  22.625 10.477  1.00 26.34 .
ATOM   2137 CG  CG  . GLU GLU A A 120 120 . -0.662  23.351 11.017  1.00 26.41 .
ATOM   2138 CD  CD  . GLU GLU A A 120 120 . 0.187   22.473 11.936  1.00 37.34 .
ATOM   2139 OE1 OE1 . GLU GLU A A 120 120 . 1.036   21.709 11.424  1.00 34.01 .
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ATOM   2140 OE2 OE2 . GLU GLU A A 120 120 . 0.002   22.537 13.170  1.00 34.00 .
ATOM   2141 N   N   . VAL VAL A A 121 121 . -0.760  24.217 7.902   1.00 28.68 .
ATOM   2142 CA  CA  . VAL VAL A A 121 121 . 0.268   24.247 6.874   1.00 18.26 .
ATOM   2143 C   C   . VAL VAL A A 121 121 . 1.545   24.589 7.631   1.00 15.73 .
ATOM   2144 O   O   . VAL VAL A A 121 121 . 1.594   25.571 8.356   1.00 22.39 .
ATOM   2145 CB  CB  . VAL VAL A A 121 121 . -0.030  25.307 5.796   1.00 22.34 .
ATOM   2146 CG1 CG1 . VAL VAL A A 121 121 . 1.180   25.476 4.887   1.00 16.48 .
ATOM   2147 CG2 CG2 . VAL VAL A A 121 121 . -1.234  24.868 4.967   1.00 26.65 .
ATOM   2148 N   N   . SER SER A A 122 122 . 2.577   23.775 7.469   1.00 17.73 .
ATOM   2149 CA  CA  . SER SER A A 122 122 . 3.819   23.998 8.198   1.00 16.30 .
ATOM   2150 C   C   . SER SER A A 122 122 . 5.046   23.909 7.313   1.00 19.39 .
ATOM   2151 O   O   . SER SER A A 122 122 . 4.944   23.637 6.117   1.00 18.32 .
ATOM   2152 CB  CB  . SER SER A A 122 122 . 3.926   22.959 9.317   1.00 19.03 .
ATOM   2153 OG  OG  . SER SER A A 122 122 . 3.913   21.644 8.768   1.00 21.56 .
ATOM   2154 N   N   . ALA ALA A A 123 123 . 6.206   24.166 7.912   1.00 16.70 .
ATOM   2155 CA  CA  . ALA ALA A A 123 123 . 7.491   24.070 7.228   1.00 16.07 .
ATOM   2156 C   C   . ALA ALA A A 123 123 . 8.106   22.862 7.935   1.00 26.23 .
ATOM   2157 O   O   . ALA ALA A A 123 123 . 8.424   22.925 9.125   1.00 20.74 .
ATOM   2158 CB  CB  . ALA ALA A A 123 123 . 8.319   25.312 7.470   1.00 20.94 .
ATOM   2159 N   N   . PHE PHE A A 124 124 . 8.253   21.762 7.206   1.00 21.78 .
ATOM   2160 CA  CA  . PHE PHE A A 124 124 . 8.747   20.532 7.802   1.00 27.82 .
ATOM   2161 C   C   . PHE PHE A A 124 124 . 10.034  19.914 7.269   1.00 21.85 .
ATOM   2162 O   O   . PHE PHE A A 124 124 . 10.472  20.181 6.149   1.00 23.70 .
ATOM   2163 CB  CB  . PHE PHE A A 124 124 . 7.610   19.489 7.760   1.00 33.17 .
ATOM   2164 CG  CG  . PHE PHE A A 124 124 . 8.069   18.065 7.962   1.00 46.24 .
ATOM   2165 CD1 CD1 . PHE PHE A A 124 124 . 8.414   17.266 6.870   1.00 50.24 .
ATOM   2166 CD2 CD2 . PHE PHE A A 124 124 . 8.184   17.534 9.242   1.00 44.44 .
ATOM   2167 CE1 CE1 . PHE PHE A A 124 124 . 8.867   15.956 7.053   1.00 55.94 .
ATOM   2168 CE2 CE2 . PHE PHE A A 124 124 . 8.636   16.229 9.438   1.00 50.06 .
ATOM   2169 CZ  CZ  . PHE PHE A A 124 124 . 8.979   15.438 8.340   1.00 50.60 .
ATOM   2170 N   N   . GLY GLY A A 125 125 . 10.644  19.095 8.126   1.00 28.18 .
ATOM   2171 CA  CA  . GLY GLY A A 125 125 . 11.848  18.361 7.784   1.00 18.74 .
ATOM   2172 C   C   . GLY GLY A A 125 125 . 13.020  19.002 7.077   1.00 19.21 .
ATOM   2173 O   O   . GLY GLY A A 125 125 . 13.539  20.031 7.508   1.00 25.00 .
ATOM   2174 N   N   . SER SER A A 126 126 . 13.442  18.355 5.992   1.00 15.81 .
ATOM   2175 CA  CA  . SER SER A A 126 126 . 14.588  18.768 5.186   1.00 17.95 .
ATOM   2176 C   C   . SER SER A A 126 126 . 14.617  17.850 3.971   1.00 11.07 .
ATOM   2177 O   O   . SER SER A A 126 126 . 13.721  17.039 3.796   1.00 23.18 .
ATOM   2178 CB  CB  . SER SER A A 126 126 . 15.879  18.560 5.977   1.00 36.05 .
ATOM   2179 OG  OG  . SER SER A A 126 126 . 16.071  17.171 6.227   1.00 32.55 .
ATOM   2180 N   N   . GLU GLU A A 127 127 . 15.651  17.953 3.145   1.00 18.71 .
ATOM   2181 CA  CA  . GLU GLU A A 127 127 . 15.739  17.092 1.972   1.00 25.35 .
ATOM   2182 C   C   . GLU GLU A A 127 127 . 15.793  15.623 2.384   1.00 32.33 .
ATOM   2183 O   O   . GLU GLU A A 127 127 . 15.260  14.748 1.694   1.00 28.03 .
ATOM   2184 CB  CB  . GLU GLU A A 127 127 . 16.974  17.445 1.139   1.00 26.06 .
ATOM   2185 CG  CG  . GLU GLU A A 127 127 . 16.824  18.743 0.372   1.00 30.24 .
ATOM   2186 CD  CD  . GLU GLU A A 127 127 . 18.035  19.083 -0.462  1.00 32.52 .
ATOM   2187 OE1 OE1 . GLU GLU A A 127 127 . 18.489  18.227 -1.254  1.00 36.10 .
ATOM   2188 OE2 OE2 . GLU GLU A A 127 127 . 18.527  20.219 -0.333  1.00 30.08 .
ATOM   2189 N   N   . ALA ALA A A 128 128 . 16.435  15.362 3.518   1.00 34.35 .
ATOM   2190 CA  CA  . ALA ALA A A 128 128 . 16.571  14.004 4.028   1.00 31.82 .
ATOM   2191 C   C   . ALA ALA A A 128 128 . 15.270  13.459 4.605   1.00 23.86 .
ATOM   2192 O   O   . ALA ALA A A 128 128 . 14.964  12.287 4.424   1.00 24.55 .
ATOM   2193 CB  CB  . ALA ALA A A 128 128 . 17.669  13.953 5.093   1.00 29.10 .
ATOM   2194 N   N   . GLU GLU A A 129 129 . 14.506  14.301 5.294   1.00 21.94 .
ATOM   2195 CA  CA  . GLU GLU A A 129 129 . 13.254  13.857 5.911   1.00 23.13 .
ATOM   2196 C   C   . GLU GLU A A 129 129 . 12.007  14.032 5.043   1.00 21.98 .
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ATOM   2197 O   O   . GLU GLU A A 129 129 . 10.924  13.589 5.417   1.00 24.31 .
ATOM   2198 CB  CB  . GLU GLU A A 129 129 . 13.005  14.615 7.214   1.00 16.36 .
ATOM   2199 CG  CG  . GLU GLU A A 129 129 . 14.116  14.538 8.242   1.00 20.95 .
ATOM   2200 CD  CD  . GLU GLU A A 129 129 . 13.968  15.613 9.293   1.00 14.22 .
ATOM   2201 OE1 OE1 . GLU GLU A A 129 129 . 13.264  15.391 10.298  1.00 30.84 .
ATOM   2202 OE2 OE2 . GLU GLU A A 129 129 . 14.541  16.704 9.097   1.00 30.97 .
ATOM   2203 N   N   . SER SER A A 130 130 . 12.153  14.704 3.911   1.00 22.32 .
ATOM   2204 CA  CA  . SER SER A A 130 130 . 11.019  14.970 3.029   1.00 27.80 .
ATOM   2205 C   C   . SER SER A A 130 130 . 10.224  13.723 2.638   1.00 28.32 .
ATOM   2206 O   O   . SER SER A A 130 130 . 10.803  12.685 2.311   1.00 26.71 .
ATOM   2207 CB  CB  . SER SER A A 130 130 . 11.511  15.679 1.766   1.00 31.02 .
ATOM   2208 OG  OG  . SER SER A A 130 130 . 10.430  16.019 0.911   1.00 27.26 .
ATOM   2209 N   N   . ASP ASP A A 131 131 . 8.897   13.838 2.682   1.00 27.76 .
ATOM   2210 CA  CA  . ASP ASP A A 131 131 . 7.997   12.744 2.305   1.00 33.60 .
ATOM   2211 C   C   . ASP ASP A A 131 131 . 6.868   13.257 1.406   1.00 28.61 .
ATOM   2212 O   O   . ASP ASP A A 131 131 . 6.863   14.429 1.025   1.00 28.09 .
ATOM   2213 CB  CB  . ASP ASP A A 131 131 . 7.409   12.053 3.547   1.00 27.91 .
ATOM   2214 CG  CG  . ASP ASP A A 131 131 . 6.581   12.987 4.416   1.00 30.93 .
ATOM   2215 OD1 OD1 . ASP ASP A A 131 131 . 5.915   13.897 3.879   1.00 32.83 .
ATOM   2216 OD2 OD2 . ASP ASP A A 131 131 . 6.576   12.793 5.651   1.00 35.70 .
ATOM   2217 N   N   . THR THR A A 132 132 . 5.912   12.389 1.076   1.00 26.65 .
ATOM   2218 CA  CA  . THR THR A A 132 132 . 4.792   12.765 0.202   1.00 25.17 .
ATOM   2219 C   C   . THR THR A A 132 132 . 3.985   13.963 0.716   1.00 23.58 .
ATOM   2220 O   O   . THR THR A A 132 132 . 3.286   14.632 -0.053  1.00 20.37 .
ATOM   2221 CB  CB  . THR THR A A 132 132 . 3.815   11.574 -0.018  1.00 21.23 .
ATOM   2222 OG1 OG1 . THR THR A A 132 132 . 3.251   11.168 1.236   1.00 25.42 .
ATOM   2223 CG2 CG2 . THR THR A A 132 132 . 4.538   10.402 -0.646  1.00 31.36 .
ATOM   2224 N   N   . GLY GLY A A 133 133 . 4.073   14.225 2.015   1.00 19.33 .
ATOM   2225 CA  CA  . GLY GLY A A 133 133 . 3.352   15.350 2.587   1.00 16.23 .
ATOM   2226 C   C   . GLY GLY A A 133 133 . 3.873   16.700 2.108   1.00 17.50 .
ATOM   2227 O   O   . GLY GLY A A 133 133 . 3.232   17.727 2.341   1.00 21.71 .
ATOM   2228 N   N   . ASP ASP A A 134 134 . 5.028   16.704 1.446   1.00 14.38 .
ATOM   2229 CA  CA  . ASP ASP A A 134 134 . 5.629   17.937 0.943   1.00 14.69 .
ATOM   2230 C   C   . ASP ASP A A 134 134 . 5.219   18.256 -0.499  1.00 22.75 .
ATOM   2231 O   O   . ASP ASP A A 134 134 . 5.552   19.331 -1.006  1.00 17.00 .
ATOM   2232 CB  CB  . ASP ASP A A 134 134 . 7.166   17.865 0.992   1.00 11.31 .
ATOM   2233 CG  CG  . ASP ASP A A 134 134 . 7.715   17.735 2.412   1.00 20.68 .
ATOM   2234 OD1 OD1 . ASP ASP A A 134 134 . 7.095   18.274 3.349   1.00 20.93 .
ATOM   2235 OD2 OD2 . ASP ASP A A 134 134 . 8.787   17.108 2.578   1.00 19.47 .
ATOM   2236 N   N   . ASP ASP A A 135 135 . 4.506   17.327 -1.148  1.00 16.29 .
ATOM   2237 CA  CA  . ASP ASP A A 135 135 . 4.085   17.507 -2.539  1.00 15.04 .
ATOM   2238 C   C   . ASP ASP A A 135 135 . 2.864   18.407 -2.746  1.00 11.53 .
ATOM   2239 O   O   . ASP ASP A A 135 135 . 1.827   18.226 -2.110  1.00 18.19 .
ATOM   2240 CB  CB  . ASP ASP A A 135 135 . 3.791   16.150 -3.203  1.00 21.95 .
ATOM   2241 CG  CG  . ASP ASP A A 135 135 . 5.044   15.313 -3.441  1.00 30.84 .
ATOM   2242 OD1 OD1 . ASP ASP A A 135 135 . 6.156   15.881 -3.552  1.00 24.55 .
ATOM   2243 OD2 OD2 . ASP ASP A A 135 135 . 4.905   14.071 -3.538  1.00 28.43 .
ATOM   2244 N   N   . TRP TRP A A 136 136 . 2.996   19.359 -3.665  1.00 16.58 .
ATOM   2245 CA  CA  . TRP TRP A A 136 136 . 1.916   20.281 -4.001  1.00 16.95 .
ATOM   2246 C   C   . TRP TRP A A 136 136 . 1.716   20.304 -5.517  1.00 23.66 .
ATOM   2247 O   O   . TRP TRP A A 136 136 . 2.683   20.415 -6.273  1.00 20.27 .
ATOM   2248 CB  CB  . TRP TRP A A 136 136 . 2.237   21.703 -3.534  1.00 17.98 .
ATOM   2249 CG  CG  . TRP TRP A A 136 136 . 2.406   21.835 -2.052  1.00 15.86 .
ATOM   2250 CD1 CD1 . TRP TRP A A 136 136 . 3.524   21.541 -1.329  1.00 20.12 .
ATOM   2251 CD2 CD2 . TRP TRP A A 136 136 . 1.426   22.288 -1.110  1.00 17.06 .
ATOM   2252 NE1 NE1 . TRP TRP A A 136 136 . 3.306   21.785 0.004   1.00 15.44 .
ATOM   2253 CE2 CE2 . TRP TRP A A 136 136 . 2.027   22.242 0.170   1.00 13.01 .
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ATOM   2254 CE3 CE3 . TRP TRP A A 136 136 . 0.103   22.727 -1.220  1.00 20.42 .
ATOM   2255 CZ2 CZ2 . TRP TRP A A 136 136 . 1.349   22.618 1.331   1.00 15.09 .
ATOM   2256 CZ3 CZ3 . TRP TRP A A 136 136 . -0.575  23.105 -0.060  1.00 25.06 .
ATOM   2257 CH2 CH2 . TRP TRP A A 136 136 . 0.052   23.046 1.199   1.00 21.57 .
ATOM   2258 N   N   . THR THR A A 137 137 . 0.465   20.196 -5.952  1.00 17.03 .
ATOM   2259 CA  CA  . THR THR A A 137 137 . 0.153   20.224 -7.373  1.00 15.09 .
ATOM   2260 C   C   . THR THR A A 137 137 . 0.025   21.671 -7.803  1.00 14.33 .
ATOM   2261 O   O   . THR THR A A 137 137 . -0.688  22.451 -7.169  1.00 21.35 .
ATOM   2262 CB  CB  . THR THR A A 137 137 . -1.174  19.508 -7.670  1.00 17.86 .
ATOM   2263 OG1 OG1 . THR THR A A 137 137 . -1.087  18.147 -7.243  1.00 18.87 .
ATOM   2264 CG2 CG2 . THR THR A A 137 137 . -1.473  19.540 -9.166  1.00 25.03 .
ATOM   2265 N   N   . VAL VAL A A 138 138 . 0.725   22.042 -8.864  1.00 14.76 .
ATOM   2266 CA  CA  . VAL VAL A A 138 138 . 0.655   23.410 -9.348  1.00 17.24 .
ATOM   2267 C   C   . VAL VAL A A 138 138 . -0.570  23.541 -10.243 1.00 22.50 .
ATOM   2268 O   O   . VAL VAL A A 138 138 . -0.708  22.825 -11.235 1.00 19.89 .
ATOM   2269 CB  CB  . VAL VAL A A 138 138 . 1.903   23.800 -10.150 1.00 15.87 .
ATOM   2270 CG1 CG1 . VAL VAL A A 138 138 . 1.859   25.281 -10.485 1.00 16.41 .
ATOM   2271 CG2 CG2 . VAL VAL A A 138 138 . 3.158   23.482 -9.341  1.00 24.07 .
ATOM   2272 N   N   . ILE ILE A A 139 139 . -1.465  24.448 -9.877  1.00 20.60 .
ATOM   2273 CA  CA  . ILE ILE A A 139 139 . -2.677  24.643 -10.655 1.00 25.88 .
ATOM   2274 C   C   . ILE ILE A A 139 139 . -2.619  25.920 -11.477 1.00 25.97 .
ATOM   2275 O   O   . ILE ILE A A 139 139 . -2.970  26.999 -10.999 1.00 28.06 .
ATOM   2276 CB  CB  . ILE ILE A A 139 139 . -3.925  24.676 -9.745  1.00 23.88 .
ATOM   2277 CG1 CG1 . ILE ILE A A 139 139 . -4.019  23.374 -8.958  1.00 25.32 .
ATOM   2278 CG2 CG2 . ILE ILE A A 139 139 . -5.193  24.853 -10.584 1.00 28.00 .
ATOM   2279 CD1 CD1 . ILE ILE A A 139 139 . -5.173  23.329 -8.000  1.00 34.89 .
ATOM   2280 N   N   . CYS CYS A A 140 140 . -2.128  25.795 -12.704 1.00 24.56 .
ATOM   2281 CA  CA  . CYS CYS A A 140 140 . -2.070  26.922 -13.615 1.00 26.01 .
ATOM   2282 C   C   . CYS CYS A A 140 140 . -2.735  26.530 -14.936 1.00 32.54 .
ATOM   2283 O   O   . CYS CYS A A 140 140 . -2.955  25.347 -15.196 1.00 30.16 .
ATOM   2284 CB  CB  . CYS CYS A A 140 140 . -0.632  27.418 -13.844 1.00 26.38 .
ATOM   2285 SG  SG  . CYS CYS A A 140 140 . 0.782   26.296 -14.153 1.00 36.60 .
ATOM   2286 N   N   . ASN ASN A A 141 141 . -3.085  27.519 -15.751 1.00 35.88 .
ATOM   2287 CA  CA  . ASN ASN A A 141 141 . -3.741  27.237 -17.023 1.00 46.03 .
ATOM   2288 C   C   . ASN ASN A A 141 141 . -2.737  27.019 -18.137 1.00 46.65 .
ATOM   2289 O   O   . ASN ASN A A 141 141 . -1.637  27.570 -18.113 1.00 40.67 .
ATOM   2290 CB  CB  . ASN ASN A A 141 141 . -4.706  28.362 -17.388 1.00 48.52 .
ATOM   2291 CG  CG  . ASN ASN A A 141 141 . -5.932  28.363 -16.513 1.00 45.87 .
ATOM   2292 OD1 OD1 . ASN ASN A A 141 141 . -6.591  27.334 -16.354 1.00 40.76 .
ATOM   2293 ND2 ND2 . ASN ASN A A 141 141 . -6.253  29.516 -15.936 1.00 48.83 .
ATOM   2294 N   N   . GLY GLY A A 142 142 . -3.131  26.218 -19.121 1.00 51.93 .
ATOM   2295 CA  CA  . GLY GLY A A 142 142 . -2.227  25.900 -20.208 1.00 57.24 .
ATOM   2296 C   C   . GLY GLY A A 142 142 . -1.367  24.789 -19.643 1.00 57.57 .
ATOM   2297 O   O   . GLY GLY A A 142 142 . -1.832  24.031 -18.791 1.00 51.15 .
ATOM   2298 N   N   . ASP ASP A A 143 143 . -0.126  24.666 -20.091 1.00 57.49 .
ATOM   2299 CA  CA  . ASP ASP A A 143 143 . 0.707   23.618 -19.528 1.00 67.33 .
ATOM   2300 C   C   . ASP ASP A A 143 143 . 2.173   23.994 -19.465 1.00 58.62 .
ATOM   2301 O   O   . ASP ASP A A 143 143 . 3.022   23.389 -20.112 1.00 70.78 .
ATOM   2302 CB  CB  . ASP ASP A A 143 143 . 0.532   22.304 -20.292 1.00 77.12 .
ATOM   2303 CG  CG  . ASP ASP A A 143 143 . 0.886   21.096 -19.437 1.00 84.59 .
ATOM   2304 OD1 OD1 . ASP ASP A A 143 143 . 2.037   21.037 -18.959 1.00 85.03 .
ATOM   2305 OD2 OD2 . ASP ASP A A 143 143 . 0.016   20.218 -19.232 1.00 89.37 .
ATOM   2306 N   N   . GLU GLU A A 144 144 . 2.449   25.018 -18.667 1.00 54.21 .
ATOM   2307 CA  CA  . GLU GLU A A 144 144 . 3.792   25.516 -18.445 1.00 52.06 .
ATOM   2308 C   C   . GLU GLU A A 144 144 . 3.672   26.743 -17.552 1.00 45.56 .
ATOM   2309 O   O   . GLU GLU A A 144 144 . 3.290   27.826 -18.000 1.00 34.20 .
ATOM   2310 CB  CB  . GLU GLU A A 144 144 . 4.480   25.863 -19.769 1.00 54.27 .
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ATOM   2311 CG  CG  . GLU GLU A A 144 144 . 5.918   25.370 -19.813 1.00 55.72 .
ATOM   2312 CD  CD  . GLU GLU A A 144 144 . 6.499   25.332 -21.212 1.00 58.85 .
ATOM   2313 OE1 OE1 . GLU GLU A A 144 144 . 7.609   24.782 -21.362 1.00 59.37 .
ATOM   2314 OE2 OE2 . GLU GLU A A 144 144 . 5.857   25.846 -22.156 1.00 59.38 .
ATOM   2315 N   N   . TRP TRP A A 145 145 . 3.971   26.533 -16.275 1.00 33.59 .
ATOM   2316 CA  CA  . TRP TRP A A 145 145 . 3.914   27.565 -15.255 1.00 28.16 .
ATOM   2317 C   C   . TRP TRP A A 145 145 . 4.623   28.829 -15.711 1.00 30.22 .
ATOM   2318 O   O   . TRP TRP A A 145 145 . 5.845   28.851 -15.818 1.00 32.16 .
ATOM   2319 CB  CB  . TRP TRP A A 145 145 . 4.562   27.031 -13.978 1.00 35.04 .
ATOM   2320 CG  CG  . TRP TRP A A 145 145 . 4.308   27.836 -12.748 1.00 32.61 .
ATOM   2321 CD1 CD1 . TRP TRP A A 145 145 . 3.571   28.982 -12.645 1.00 36.35 .
ATOM   2322 CD2 CD2 . TRP TRP A A 145 145 . 4.788   27.546 -11.433 1.00 29.98 .
ATOM   2323 NE1 NE1 . TRP TRP A A 145 145 . 3.565   29.423 -11.340 1.00 36.35 .
ATOM   2324 CE2 CE2 . TRP TRP A A 145 145 . 4.304   28.560 -10.574 1.00 29.32 .
ATOM   2325 CE3 CE3 . TRP TRP A A 145 145 . 5.584   26.525 -10.895 1.00 26.81 .
ATOM   2326 CZ2 CZ2 . TRP TRP A A 145 145 . 4.591   28.581 -9.206  1.00 25.27 .
ATOM   2327 CZ3 CZ3 . TRP TRP A A 145 145 . 5.869   26.549 -9.533  1.00 18.38 .
ATOM   2328 CH2 CH2 . TRP TRP A A 145 145 . 5.372   27.572 -8.705  1.00 25.18 .
ATOM   2329 N   N   . LEU LEU A A 146 146 . 3.847   29.878 -15.979 1.00 28.93 .
ATOM   2330 CA  CA  . LEU LEU A A 146 146 . 4.395   31.158 -16.422 1.00 33.55 .
ATOM   2331 C   C   . LEU LEU A A 146 146 . 4.501   32.099 -15.224 1.00 37.58 .
ATOM   2332 O   O   . LEU LEU A A 146 146 . 3.601   32.164 -14.384 1.00 38.21 .
ATOM   2333 CB  CB  . LEU LEU A A 146 146 . 3.489   31.785 -17.487 1.00 36.54 .
ATOM   2334 CG  CG  . LEU LEU A A 146 146 . 3.125   30.948 -18.718 1.00 39.91 .
ATOM   2335 CD1 CD1 . LEU LEU A A 146 146 . 2.103   31.704 -19.558 1.00 37.99 .
ATOM   2336 CD2 CD2 . LEU LEU A A 146 146 . 4.371   30.638 -19.535 1.00 38.87 .
ATOM   2337 N   N   . GLU GLU A A 147 147 . 5.605   32.830 -15.151 1.00 35.71 .
ATOM   2338 CA  CA  . GLU GLU A A 147 147 . 5.842   33.755 -14.054 1.00 33.26 .
ATOM   2339 C   C   . GLU GLU A A 147 147 . 4.837   34.905 -14.015 1.00 40.18 .
ATOM   2340 O   O   . GLU GLU A A 147 147 . 4.712   35.594 -13.002 1.00 31.21 .
ATOM   2341 CB  CB  . GLU GLU A A 147 147 . 7.261   34.313 -14.165 1.00 43.38 .
ATOM   2342 CG  CG  . GLU GLU A A 147 147 . 7.690   35.179 -12.997 1.00 54.33 .
ATOM   2343 CD  CD  . GLU GLU A A 147 147 . 9.182   35.441 -13.002 1.00 57.84 .
ATOM   2344 OE1 OE1 . GLU GLU A A 147 147 . 9.944   34.458 -13.128 1.00 61.08 .
ATOM   2345 OE2 OE2 . GLU GLU A A 147 147 . 9.589   36.616 -12.873 1.00 59.41 .
ATOM   2346 N   N   . SER SER A A 148 148 . 4.112   35.093 -15.114 1.00 38.98 .
ATOM   2347 CA  CA  . SER SER A A 148 148 . 3.141   36.180 -15.228 1.00 42.25 .
ATOM   2348 C   C   . SER SER A A 148 148 . 1.700   35.786 -14.925 1.00 43.42 .
ATOM   2349 O   O   . SER SER A A 148 148 . 0.789   36.604 -15.057 1.00 45.87 .
ATOM   2350 CB  CB  . SER SER A A 148 148 . 3.191   36.756 -16.639 1.00 46.08 .
ATOM   2351 OG  OG  . SER SER A A 148 148 . 2.737   35.789 -17.572 1.00 45.05 .
ATOM   2352 N   N   . GLU GLU A A 149 149 . 1.477   34.546 -14.522 1.00 37.15 .
ATOM   2353 CA  CA  . GLU GLU A A 149 149 . 0.115   34.128 -14.257 1.00 35.84 .
ATOM   2354 C   C   . GLU GLU A A 149 149 . -0.147  33.851 -12.790 1.00 35.39 .
ATOM   2355 O   O   . GLU GLU A A 149 149 . 0.781   33.659 -12.001 1.00 30.63 .
ATOM   2356 CB  CB  . GLU GLU A A 149 149 . -0.216  32.886 -15.083 1.00 39.34 .
ATOM   2357 CG  CG  . GLU GLU A A 149 149 . 0.354   31.597 -14.528 1.00 54.82 .
ATOM   2358 CD  CD  . GLU GLU A A 149 149 . 0.320   30.471 -15.541 1.00 63.46 .
ATOM   2359 OE1 OE1 . GLU GLU A A 149 149 . 1.243   30.406 -16.379 1.00 60.83 .
ATOM   2360 OE2 OE2 . GLU GLU A A 149 149 . -0.633  29.661 -15.509 1.00 64.54 .
ATOM   2361 N   N   . GLN GLN A A 150 150 . -1.422  33.854 -12.424 1.00 31.44 .
ATOM   2362 CA  CA  . GLN GLN A A 150 150 . -1.788  33.561 -11.058 1.00 30.28 .
ATOM   2363 C   C   . GLN GLN A A 150 150 . -1.933  32.053 -11.039 1.00 32.33 .
ATOM   2364 O   O   . GLN GLN A A 150 150 . -2.091  31.429 -12.090 1.00 33.78 .
ATOM   2365 CB  CB  . GLN GLN A A 150 150 . -3.108  34.231 -10.677 1.00 30.18 .
ATOM   2366 CG  CG  . GLN GLN A A 150 150 . -3.093  35.736 -10.829 1.00 28.97 .
ATOM   2367 CD  CD  . GLN GLN A A 150 150 . -4.005  36.435 -9.845  1.00 22.84 .
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ATOM   2368 OE1 OE1 . GLN GLN A A 150 150 . -5.049  35.911 -9.465  1.00 30.20 .
ATOM   2369 NE2 NE2 . GLN GLN A A 150 150 . -3.617  37.635 -9.432  1.00 31.81 .
ATOM   2370 N   N   . PHE PHE A A 151 151 . -1.875  31.466 -9.853  1.00 20.93 .
ATOM   2371 CA  CA  . PHE PHE A A 151 151 . -1.982  30.028 -9.736  1.00 24.48 .
ATOM   2372 C   C   . PHE PHE A A 151 151 . -2.422  29.640 -8.341  1.00 24.54 .
ATOM   2373 O   O   . PHE PHE A A 151 151 . -2.466  30.472 -7.436  1.00 23.90 .
ATOM   2374 CB  CB  . PHE PHE A A 151 151 . -0.618  29.380 -10.024 1.00 24.29 .
ATOM   2375 CG  CG  . PHE PHE A A 151 151 . 0.468   29.775 -9.043  1.00 21.96 .
ATOM   2376 CD1 CD1 . PHE PHE A A 151 151 . 1.170   30.970 -9.195  1.00 31.49 .
ATOM   2377 CD2 CD2 . PHE PHE A A 151 151 . 0.765   28.965 -7.949  1.00 21.17 .
ATOM   2378 CE1 CE1 . PHE PHE A A 151 151 . 2.155   31.355 -8.265  1.00 23.83 .
ATOM   2379 CE2 CE2 . PHE PHE A A 151 151 . 1.742   29.341 -7.019  1.00 17.41 .
ATOM   2380 CZ  CZ  . PHE PHE A A 151 151 . 2.435   30.534 -7.176  1.00 16.57 .
ATOM   2381 N   N   . LYS LYS A A 152 152 . -2.769  28.372 -8.180  1.00 20.08 .
ATOM   2382 CA  CA  . LYS LYS A A 152 152 . -3.138  27.853 -6.880  1.00 21.94 .
ATOM   2383 C   C   . LYS LYS A A 152 152 . -2.202  26.677 -6.635  1.00 21.91 .
ATOM   2384 O   O   . LYS LYS A A 152 152 . -1.544  26.188 -7.563  1.00 22.08 .
ATOM   2385 CB  CB  . LYS LYS A A 152 152 . -4.594  27.382 -6.865  1.00 33.16 .
ATOM   2386 CG  CG  . LYS LYS A A 152 152 . -5.603  28.512 -6.930  1.00 25.45 .
ATOM   2387 CD  CD  . LYS LYS A A 152 152 . -7.018  27.989 -7.035  1.00 41.56 .
ATOM   2388 CE  CE  . LYS LYS A A 152 152 . -7.474  27.896 -8.488  1.00 47.49 .
ATOM   2389 NZ  NZ  . LYS LYS A A 152 152 . -7.620  29.243 -9.118  1.00 48.62 .
ATOM   2390 N   N   . LEU LEU A A 153 153 . -2.120  26.241 -5.388  1.00 15.13 .
ATOM   2391 CA  CA  . LEU LEU A A 153 153 . -1.288  25.105 -5.032  1.00 17.30 .
ATOM   2392 C   C   . LEU LEU A A 153 153 . -2.165  24.176 -4.230  1.00 13.83 .
ATOM   2393 O   O   . LEU LEU A A 153 153 . -2.640  24.529 -3.152  1.00 22.19 .
ATOM   2394 CB  CB  . LEU LEU A A 153 153 . -0.082  25.531 -4.191  1.00 16.98 .
ATOM   2395 CG  CG  . LEU LEU A A 153 153 . 1.087   26.212 -4.902  1.00 20.49 .
ATOM   2396 CD1 CD1 . LEU LEU A A 153 153 . 2.138   26.596 -3.855  1.00 20.54 .
ATOM   2397 CD2 CD2 . LEU LEU A A 153 153 . 1.698   25.281 -5.945  1.00 19.18 .
ATOM   2398 N   N   . ARG ARG A A 154 154 . -2.378  22.981 -4.759  1.00 14.51 .
ATOM   2399 CA  CA  . ARG ARG A A 154 154 . -3.224  22.016 -4.085  1.00 18.83 .
ATOM   2400 C   C   . ARG ARG A A 154 154 . -2.398  20.890 -3.482  1.00 17.50 .
ATOM   2401 O   O   . ARG ARG A A 154 154 . -1.681  20.192 -4.195  1.00 18.72 .
ATOM   2402 CB  CB  . ARG ARG A A 154 154 . -4.237  21.452 -5.085  1.00 21.15 .
ATOM   2403 CG  CG  . ARG ARG A A 154 154 . -5.421  20.768 -4.448  1.00 23.17 .
ATOM   2404 CD  CD  . ARG ARG A A 154 154 . -6.440  20.327 -5.512  1.00 24.51 .
ATOM   2405 NE  NE  . ARG ARG A A 154 154 . -7.684  19.871 -4.901  1.00 26.09 .
ATOM   2406 CZ  CZ  . ARG ARG A A 154 154 . -8.759  19.481 -5.583  1.00 30.30 .
ATOM   2407 NH1 NH1 . ARG ARG A A 154 154 . -8.749  19.485 -6.910  1.00 21.54 .
ATOM   2408 NH2 NH2 . ARG ARG A A 154 154 . -9.847  19.098 -4.936  1.00 24.59 .
ATOM   2409 N   N   . HIS HIS A A 155 155 . -2.501  20.714 -2.168  1.00 15.47 .
ATOM   2410 CA  CA  . HIS HIS A A 155 155 . -1.750  19.666 -1.489  1.00 17.47 .
ATOM   2411 C   C   . HIS HIS A A 155 155 . -2.115  18.323 -2.114  1.00 29.13 .
ATOM   2412 O   O   . HIS HIS A A 155 155 . -3.286  17.934 -2.140  1.00 23.05 .
ATOM   2413 CB  CB  . HIS HIS A A 155 155 . -2.077  19.652 -0.001  1.00 20.13 .
ATOM   2414 CG  CG  . HIS HIS A A 155 155 . -1.115  18.851 0.819   1.00 28.32 .
ATOM   2415 ND1 ND1 . HIS HIS A A 155 155 . -1.522  17.867 1.696   1.00 26.70 .
ATOM   2416 CD2 CD2 . HIS HIS A A 155 155 . 0.235   18.909 0.918   1.00 18.85 .
ATOM   2417 CE1 CE1 . HIS HIS A A 155 155 . -0.464  17.356 2.302   1.00 26.66 .
ATOM   2418 NE2 NE2 . HIS HIS A A 155 155 . 0.614   17.971 1.848   1.00 24.06 .
ATOM   2419 N   N   . ALA ALA A A 156 156 . -1.105  17.620 -2.610  1.00 20.80 .
ATOM   2420 CA  CA  . ALA ALA A A 156 156 . -1.314  16.341 -3.263  1.00 26.82 .
ATOM   2421 C   C   . ALA ALA A A 156 156 . -1.967  15.272 -2.392  1.00 31.22 .
ATOM   2422 O   O   . ALA ALA A A 156 156 . -2.929  14.630 -2.813  1.00 26.75 .
ATOM   2423 CB  CB  . ALA ALA A A 156 156 . 0.005   15.825 -3.807  1.00 22.75 .
ATOM   2424 N   N   . VAL VAL A A 157 157 . -1.451  15.082 -1.183  1.00 24.58 .
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ATOM   2425 CA  CA  . VAL VAL A A 157 157 . -1.978  14.057 -0.296  1.00 28.01 .
ATOM   2426 C   C   . VAL VAL A A 157 157 . -3.366  14.316 0.283   1.00 35.31 .
ATOM   2427 O   O   . VAL VAL A A 157 157 . -4.182  13.398 0.351   1.00 28.08 .
ATOM   2428 CB  CB  . VAL VAL A A 157 157 . -0.990  13.774 0.852   1.00 24.27 .
ATOM   2429 CG1 CG1 . VAL VAL A A 157 157 . -1.615  12.839 1.880   1.00 30.76 .
ATOM   2430 CG2 CG2 . VAL VAL A A 157 157 . 0.279   13.156 0.286   1.00 22.61 .
ATOM   2431 N   N   . THR THR A A 158 158 . -3.645  15.551 0.698   1.00 35.67 .
ATOM   2432 CA  CA  . THR THR A A 158 158 . -4.950  15.870 1.277   1.00 30.59 .
ATOM   2433 C   C   . THR THR A A 158 158 . -5.911  16.580 0.326   1.00 29.07 .
ATOM   2434 O   O   . THR THR A A 158 158 . -7.123  16.534 0.518   1.00 34.19 .
ATOM   2435 CB  CB  . THR THR A A 158 158 . -4.806  16.755 2.522   1.00 29.26 .
ATOM   2436 OG1 OG1 . THR THR A A 158 158 . -4.178  17.986 2.151   1.00 28.82 .
ATOM   2437 CG2 CG2 . THR THR A A 158 158 . -3.980  16.059 3.583   1.00 26.59 .
ATOM   2438 N   N   . GLY GLY A A 159 159 . -5.376  17.249 -0.688  1.00 29.90 .
ATOM   2439 CA  CA  . GLY GLY A A 159 159 . -6.225  17.957 -1.632  1.00 26.32 .
ATOM   2440 C   C   . GLY GLY A A 159 159 . -6.556  19.384 -1.222  1.00 26.34 .
ATOM   2441 O   O   . GLY GLY A A 159 159 . -7.235  20.107 -1.956  1.00 31.61 .
ATOM   2442 N   N   . SER SER A A 160 160 . -6.076  19.798 -0.053  1.00 28.60 .
ATOM   2443 CA  CA  . SER SER A A 160 160 . -6.341  21.144 0.443   1.00 23.42 .
ATOM   2444 C   C   . SER SER A A 160 160 . -5.551  22.180 -0.332  1.00 28.84 .
ATOM   2445 O   O   . SER SER A A 160 160 . -4.408  21.939 -0.729  1.00 23.36 .
ATOM   2446 CB  CB  . SER SER A A 160 160 . -5.987  21.247 1.927   1.00 32.10 .
ATOM   2447 OG  OG  . SER SER A A 160 160 . -6.621  20.219 2.667   1.00 31.58 .
ATOM   2448 N   N   . TYR TYR A A 161 161 . -6.171  23.336 -0.543  1.00 23.65 .
ATOM   2449 CA  CA  . TYR TYR A A 161 161 . -5.543  24.425 -1.263  1.00 17.93 .
ATOM   2450 C   C   . TYR TYR A A 161 161 . -4.767  25.308 -0.291  1.00 23.69 .
ATOM   2451 O   O   . TYR TYR A A 161 161 . -5.220  25.567 0.828   1.00 21.29 .
ATOM   2452 CB  CB  . TYR TYR A A 161 161 . -6.607  25.260 -1.964  1.00 21.44 .
ATOM   2453 CG  CG  . TYR TYR A A 161 161 . -7.327  24.533 -3.076  1.00 30.77 .
ATOM   2454 CD1 CD1 . TYR TYR A A 161 161 . -6.779  24.458 -4.353  1.00 21.42 .
ATOM   2455 CD2 CD2 . TYR TYR A A 161 161 . -8.563  23.918 -2.851  1.00 32.94 .
ATOM   2456 CE1 CE1 . TYR TYR A A 161 161 . -7.440  23.793 -5.384  1.00 23.20 .
ATOM   2457 CE2 CE2 . TYR TYR A A 161 161 . -9.234  23.247 -3.878  1.00 31.54 .
ATOM   2458 CZ  CZ  . TYR TYR A A 161 161 . -8.665  23.189 -5.141  1.00 28.73 .
ATOM   2459 OH  OH  . TYR TYR A A 161 161 . -9.312  22.521 -6.165  1.00 35.08 .
ATOM   2460 N   N   . LEU LEU A A 162 162 . -3.590  25.754 -0.709  1.00 19.19 .
ATOM   2461 CA  CA  . LEU LEU A A 162 162 . -2.785  26.634 0.134   1.00 18.10 .
ATOM   2462 C   C   . LEU LEU A A 162 162 . -3.559  27.948 0.116   1.00 13.12 .
ATOM   2463 O   O   . LEU LEU A A 162 162 . -3.795  28.502 -0.956  1.00 15.98 .
ATOM   2464 CB  CB  . LEU LEU A A 162 162 . -1.396  26.830 -0.486  1.00 10.84 .
ATOM   2465 CG  CG  . LEU LEU A A 162 162 . -0.439  27.768 0.253   1.00 11.79 .
ATOM   2466 CD1 CD1 . LEU LEU A A 162 162 . -0.038  27.144 1.577   1.00 13.94 .
ATOM   2467 CD2 CD2 . LEU LEU A A 162 162 . 0.786   28.018 -0.615  1.00 13.26 .
ATOM   2468 N   N   . SER SER A A 163 163 . -3.965  28.441 1.283   1.00 17.70 .
ATOM   2469 CA  CA  . SER SER A A 163 163 . -4.748  29.676 1.321   1.00 19.54 .
ATOM   2470 C   C   . SER SER A A 163 163 . -4.541  30.507 2.572   1.00 18.67 .
ATOM   2471 O   O   . SER SER A A 163 163 . -4.176  29.985 3.625   1.00 20.90 .
ATOM   2472 CB  CB  . SER SER A A 163 163 . -6.248  29.357 1.236   1.00 20.70 .
ATOM   2473 OG  OG  . SER SER A A 163 163 . -6.551  28.450 0.192   1.00 24.20 .
ATOM   2474 N   N   . LEU LEU A A 164 164 . -4.771  31.809 2.445   1.00 21.67 .
ATOM   2475 CA  CA  . LEU LEU A A 164 164 . -4.691  32.689 3.599   1.00 19.45 .
ATOM   2476 C   C   . LEU LEU A A 164 164 . -6.078  32.529 4.225   1.00 24.69 .
ATOM   2477 O   O   . LEU LEU A A 164 164 . -7.096  32.646 3.532   1.00 29.08 .
ATOM   2478 CB  CB  . LEU LEU A A 164 164 . -4.470  34.143 3.170   1.00 20.89 .
ATOM   2479 CG  CG  . LEU LEU A A 164 164 . -3.128  34.460 2.502   1.00 22.75 .
ATOM   2480 CD1 CD1 . LEU LEU A A 164 164 . -2.985  35.969 2.325   1.00 18.25 .
ATOM   2481 CD2 CD2 . LEU LEU A A 164 164 . -1.994  33.922 3.363   1.00 24.82 .
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ATOM   2482 N   N   . SER SER A A 165 165 . -6.117  32.224 5.515   1.00 23.87 .
ATOM   2483 CA  CA  . SER SER A A 165 165 . -7.381  32.035 6.222   1.00 28.68 .
ATOM   2484 C   C   . SER SER A A 165 165 . -8.107  33.357 6.399   1.00 32.54 .
ATOM   2485 O   O   . SER SER A A 165 165 . -9.321  33.386 6.584   1.00 33.83 .
ATOM   2486 CB  CB  . SER SER A A 165 165 . -7.129  31.432 7.601   1.00 22.88 .
ATOM   2487 OG  OG  . SER SER A A 165 165 . -6.500  32.384 8.436   1.00 23.79 .
ATOM   2488 N   N   . GLY GLY A A 166 166 . -7.353  34.451 6.346   1.00 34.31 .
ATOM   2489 CA  CA  . GLY GLY A A 166 166 . -7.941  35.766 6.521   1.00 34.98 .
ATOM   2490 C   C   . GLY GLY A A 166 166 . -7.682  36.246 7.936   1.00 36.96 .
ATOM   2491 O   O   . GLY GLY A A 166 166 . -7.992  37.380 8.296   1.00 37.58 .
ATOM   2492 N   N   . GLN GLN A A 167 167 . -7.111  35.363 8.747   1.00 38.43 .
ATOM   2493 CA  CA  . GLN GLN A A 167 167 . -6.784  35.690 10.124  1.00 34.20 .
ATOM   2494 C   C   . GLN GLN A A 167 167 . -5.351  36.188 10.196  1.00 33.80 .
ATOM   2495 O   O   . GLN GLN A A 167 167 . -4.536  35.887 9.325   1.00 33.53 .
ATOM   2496 CB  CB  . GLN GLN A A 167 167 . -6.936  34.454 11.013  1.00 41.70 .
ATOM   2497 CG  CG  . GLN GLN A A 167 167 . -8.373  34.025 11.223  1.00 39.98 .
ATOM   2498 CD  CD  . GLN GLN A A 167 167 . -9.194  35.120 11.870  1.00 44.06 .
ATOM   2499 OE1 OE1 . GLN GLN A A 167 167 . -8.883  35.576 12.971  1.00 41.63 .
ATOM   2500 NE2 NE2 . GLN GLN A A 167 167 . -10.243 35.556 11.185  1.00 42.38 .
ATOM   2501 N   N   . GLN GLN A A 168 168 . -5.053  36.954 11.238  1.00 32.46 .
ATOM   2502 CA  CA  . GLN GLN A A 168 168 . -3.714  37.488 11.448  1.00 28.84 .
ATOM   2503 C   C   . GLN GLN A A 168 168 . -3.239  37.090 12.840  1.00 34.11 .
ATOM   2504 O   O   . GLN GLN A A 168 168 . -4.047  36.943 13.760  1.00 36.68 .
ATOM   2505 CB  CB  . GLN GLN A A 168 168 . -3.719  39.011 11.328  1.00 30.89 .
ATOM   2506 CG  CG  . GLN GLN A A 168 168 . -3.897  39.539 9.919   1.00 28.20 .
ATOM   2507 CD  CD  . GLN GLN A A 168 168 . -3.769  41.054 9.854   1.00 40.50 .
ATOM   2508 OE1 OE1 . GLN GLN A A 168 168 . -2.911  41.643 10.514  1.00 28.64 .
ATOM   2509 NE2 NE2 . GLN GLN A A 168 168 . -4.613  41.689 9.047   1.00 36.77 .
ATOM   2510 N   N   . PHE PHE A A 169 169 . -1.932  36.917 12.992  1.00 24.96 .
ATOM   2511 CA  CA  . PHE PHE A A 169 169 . -1.363  36.546 14.277  1.00 27.15 .
ATOM   2512 C   C   . PHE PHE A A 169 169 . -1.045  37.774 15.109  1.00 33.71 .
ATOM   2513 O   O   . PHE PHE A A 169 169 . -1.000  38.897 14.598  1.00 24.98 .
ATOM   2514 CB  CB  . PHE PHE A A 169 169 . -0.065  35.771 14.092  1.00 25.98 .
ATOM   2515 CG  CG  . PHE PHE A A 169 169 . -0.246  34.401 13.527  1.00 27.91 .
ATOM   2516 CD1 CD1 . PHE PHE A A 169 169 . -0.794  33.386 14.300  1.00 28.34 .
ATOM   2517 CD2 CD2 . PHE PHE A A 169 169 . 0.146   34.118 12.223  1.00 27.44 .
ATOM   2518 CE1 CE1 . PHE PHE A A 169 169 . -0.951  32.105 13.788  1.00 22.98 .
ATOM   2519 CE2 CE2 . PHE PHE A A 169 169 . -0.006  32.835 11.697  1.00 32.23 .
ATOM   2520 CZ  CZ  . PHE PHE A A 169 169 . -0.557  31.827 12.484  1.00 34.10 .
ATOM   2521 N   N   . GLY GLY A A 170 170 . -0.821  37.534 16.398  1.00 35.76 .
ATOM   2522 CA  CA  . GLY GLY A A 170 170 . -0.449  38.591 17.316  1.00 27.14 .
ATOM   2523 C   C   . GLY GLY A A 170 170 . 1.017   38.323 17.587  1.00 34.55 .
ATOM   2524 O   O   . GLY GLY A A 170 170 . 1.647   37.575 16.835  1.00 27.88 .
ATOM   2525 N   N   . ARG ARG A A 171 171 . 1.570   38.907 18.645  1.00 32.53 .
ATOM   2526 CA  CA  . ARG ARG A A 171 171 . 2.979   38.691 18.976  1.00 29.52 .
ATOM   2527 C   C   . ARG ARG A A 171 171 . 3.225   37.192 19.147  1.00 2.79  .
ATOM   2528 O   O   . ARG ARG A A 171 171 . 2.304   36.442 19.462  1.00 33.34 .
ATOM   2529 CB  CB  . ARG ARG A A 171 171 . 3.334   39.416 20.275  1.00 33.63 .
ATOM   2530 CG  CG  . ARG ARG A A 171 171 . 2.895   38.687 21.539  1.00 41.53 .
ATOM   2531 CD  CD  . ARG ARG A A 171 171 . 1.399   38.403 21.535  1.00 57.40 .
ATOM   2532 NE  NE  . ARG ARG A A 171 171 . 0.606   39.624 21.411  1.00 65.82 .
ATOM   2533 CZ  CZ  . ARG ARG A A 171 171 . 0.687   40.655 22.249  1.00 70.52 .
ATOM   2534 NH1 NH1 . ARG ARG A A 171 171 . 1.528   40.615 23.275  1.00 65.83 .
ATOM   2535 NH2 NH2 . ARG ARG A A 171 171 . -0.080  41.722 22.068  1.00 71.20 .
ATOM   2536 N   N   . PRO PRO A A 172 172 . 4.471   36.734 18.953  1.00 26.19 .
ATOM   2537 CA  CA  . PRO PRO A A 172 172 . 5.684   37.472 18.589  1.00 26.84 .
ATOM   2538 C   C   . PRO PRO A A 172 172 . 5.826   37.712 17.086  1.00 25.46 .
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ATOM   2539 O   O   . PRO PRO A A 172 172 . 6.896   38.097 16.615  1.00 28.10 .
ATOM   2540 CB  CB  . PRO PRO A A 172 172 . 6.789   36.570 19.114  1.00 34.21 .
ATOM   2541 CG  CG  . PRO PRO A A 172 172 . 6.236   35.211 18.798  1.00 28.42 .
ATOM   2542 CD  CD  . PRO PRO A A 172 172 . 4.797   35.327 19.255  1.00 30.70 .
ATOM   2543 N   N   . ILE ILE A A 173 173 . 4.750   37.472 16.341  1.00 24.66 .
ATOM   2544 CA  CA  . ILE ILE A A 173 173 . 4.746   37.659 14.887  1.00 25.37 .
ATOM   2545 C   C   . ILE ILE A A 173 173 . 3.519   38.478 14.511  1.00 24.11 .
ATOM   2546 O   O   . ILE ILE A A 173 173 . 2.754   38.109 13.622  1.00 26.04 .
ATOM   2547 CB  CB  . ILE ILE A A 173 173 . 4.697   36.289 14.145  1.00 23.12 .
ATOM   2548 CG1 CG1 . ILE ILE A A 173 173 . 3.592   35.411 14.735  1.00 21.38 .
ATOM   2549 CG2 CG2 . ILE ILE A A 173 173 . 6.052   35.591 14.242  1.00 21.28 .
ATOM   2550 CD1 CD1 . ILE ILE A A 173 173 . 3.371   34.090 13.984  1.00 24.83 .
ATOM   2551 N   N   . HIS HIS A A 174 174 . 3.350   39.603 15.199  1.00 24.41 .
ATOM   2552 CA  CA  . HIS HIS A A 174 174 . 2.215   40.490 14.997  1.00 19.18 .
ATOM   2553 C   C   . HIS HIS A A 174 174 . 1.957   40.991 13.583  1.00 17.58 .
ATOM   2554 O   O   . HIS HIS A A 174 174 . 2.839   41.541 12.926  1.00 19.98 .
ATOM   2555 CB  CB  . HIS HIS A A 174 174 . 2.335   41.691 15.942  1.00 30.56 .
ATOM   2556 CG  CG  . HIS HIS A A 174 174 . 1.284   42.734 15.731  1.00 27.69 .
ATOM   2557 ND1 ND1 . HIS HIS A A 174 174 . 1.341   43.649 14.703  1.00 28.94 .
ATOM   2558 CD2 CD2 . HIS HIS A A 174 174 . 0.137   42.992 16.403  1.00 25.57 .
ATOM   2559 CE1 CE1 . HIS HIS A A 174 174 . 0.274   44.427 14.749  1.00 27.31 .
ATOM   2560 NE2 NE2 . HIS HIS A A 174 174 . -0.473  44.048 15.771  1.00 27.47 .
ATOM   2561 N   N   . GLY GLY A A 175 175 . 0.723   40.813 13.130  1.00 19.47 .
ATOM   2562 CA  CA  . GLY GLY A A 175 175 . 0.341   41.276 11.810  1.00 24.14 .
ATOM   2563 C   C   . GLY GLY A A 175 175 . 0.550   40.278 10.685  1.00 25.33 .
ATOM   2564 O   O   . GLY GLY A A 175 175 . 0.095   40.507 9.567   1.00 28.06 .
ATOM   2565 N   N   . GLN GLN A A 176 176 . 1.240   39.179 10.959  1.00 20.96 .
ATOM   2566 CA  CA  . GLN GLN A A 176 176 . 1.478   38.180 9.922   1.00 17.60 .
ATOM   2567 C   C   . GLN GLN A A 176 176 . 0.203   37.374 9.684   1.00 21.38 .
ATOM   2568 O   O   . GLN GLN A A 176 176 . -0.511  37.028 10.629  1.00 21.39 .
ATOM   2569 CB  CB  . GLN GLN A A 176 176 . 2.646   37.279 10.327  1.00 23.06 .
ATOM   2570 CG  CG  . GLN GLN A A 176 176 . 3.968   38.043 10.409  1.00 23.13 .
ATOM   2571 CD  CD  . GLN GLN A A 176 176 . 5.165   37.159 10.672  1.00 19.15 .
ATOM   2572 OE1 OE1 . GLN GLN A A 176 176 . 5.220   36.020 10.221  1.00 20.30 .
ATOM   2573 NE2 NE2 . GLN GLN A A 176 176 . 6.151   37.695 11.386  1.00 26.10 .
ATOM   2574 N   N   . ARG ARG A A 177 177 . -0.095  37.092 8.419   1.00 20.95 .
ATOM   2575 CA  CA  . ARG ARG A A 177 177 . -1.312  36.358 8.078   1.00 23.16 .
ATOM   2576 C   C   . ARG ARG A A 177 177 . -1.215  34.844 8.238   1.00 22.67 .
ATOM   2577 O   O   . ARG ARG A A 177 177 . -0.179  34.241 7.947   1.00 19.49 .
ATOM   2578 CB  CB  . ARG ARG A A 177 177 . -1.746  36.688 6.647   1.00 25.76 .
ATOM   2579 CG  CG  . ARG ARG A A 177 177 . -2.163  38.141 6.423   1.00 22.12 .
ATOM   2580 CD  CD  . ARG ARG A A 177 177 . -0.964  39.065 6.319   1.00 24.32 .
ATOM   2581 NE  NE  . ARG ARG A A 177 177 . -1.288  40.322 5.646   1.00 28.43 .
ATOM   2582 CZ  CZ  . ARG ARG A A 177 177 . -1.854  41.373 6.230   1.00 40.37 .
ATOM   2583 NH1 NH1 . ARG ARG A A 177 177 . -2.167  41.336 7.517   1.00 41.82 .
ATOM   2584 NH2 NH2 . ARG ARG A A 177 177 . -2.109  42.469 5.521   1.00 39.87 .
ATOM   2585 N   N   . GLU GLU A A 178 178 . -2.306  34.237 8.707   1.00 17.81 .
ATOM   2586 CA  CA  . GLU GLU A A 178 178 . -2.348  32.791 8.889   1.00 24.64 .
ATOM   2587 C   C   . GLU GLU A A 178 178 . -2.487  32.113 7.529   1.00 19.11 .
ATOM   2588 O   O   . GLU GLU A A 178 178 . -3.314  32.511 6.707   1.00 21.73 .
ATOM   2589 CB  CB  . GLU GLU A A 178 178 . -3.532  32.381 9.782   1.00 28.23 .
ATOM   2590 CG  CG  . GLU GLU A A 178 178 . -3.709  30.866 9.905   1.00 28.48 .
ATOM   2591 CD  CD  . GLU GLU A A 178 178 . -5.008  30.462 10.602  1.00 37.76 .
ATOM   2592 OE1 OE1 . GLU GLU A A 178 178 . -6.037  31.137 10.386  1.00 25.51 .
ATOM   2593 OE2 OE2 . GLU GLU A A 178 178 . -5.000  29.460 11.350  1.00 29.77 .
ATOM   2594 N   N   . VAL VAL A A 179 179 . -1.652  31.105 7.303   1.00 22.11 .
ATOM   2595 CA  CA  . VAL VAL A A 179 179 . -1.659  30.325 6.074   1.00 16.88 .
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ATOM   2596 C   C   . VAL VAL A A 179 179 . -2.212  28.945 6.448   1.00 16.94 .
ATOM   2597 O   O   . VAL VAL A A 179 179 . -1.755  28.321 7.417   1.00 21.65 .
ATOM   2598 CB  CB  . VAL VAL A A 179 179 . -0.232  30.163 5.508   1.00 20.80 .
ATOM   2599 CG1 CG1 . VAL VAL A A 179 179 . -0.250  29.202 4.323   1.00 19.35 .
ATOM   2600 CG2 CG2 . VAL VAL A A 179 179 . 0.327   31.517 5.078   1.00 14.44 .
ATOM   2601 N   N   . VAL VAL A A 180 180 . -3.192  28.467 5.694   1.00 19.97 .
ATOM   2602 CA  CA  . VAL VAL A A 180 180 . -3.794  27.175 5.998   1.00 24.10 .
ATOM   2603 C   C   . VAL VAL A A 180 180 . -4.199  26.402 4.750   1.00 26.36 .
ATOM   2604 O   O   . VAL VAL A A 180 180 . -4.016  26.864 3.623   1.00 21.67 .
ATOM   2605 CB  CB  . VAL VAL A A 180 180 . -5.065  27.330 6.860   1.00 23.18 .
ATOM   2606 CG1 CG1 . VAL VAL A A 180 180 . -4.766  28.140 8.114   1.00 25.66 .
ATOM   2607 CG2 CG2 . VAL VAL A A 180 180 . -6.162  27.989 6.041   1.00 23.75 .
ATOM   2608 N   N   . GLY GLY A A 181 181 . -4.769  25.222 4.980   1.00 26.58 .
ATOM   2609 CA  CA  . GLY GLY A A 181 181 . -5.219  24.377 3.892   1.00 26.55 .
ATOM   2610 C   C   . GLY GLY A A 181 181 . -6.735  24.355 3.905   1.00 24.92 .
ATOM   2611 O   O   . GLY GLY A A 181 181 . -7.345  23.923 4.880   1.00 29.10 .
ATOM   2612 N   N   . THR THR A A 182 182 . -7.345  24.838 2.832   1.00 24.48 .
ATOM   2613 CA  CA  . THR THR A A 182 182 . -8.794  24.868 2.754   1.00 30.93 .
ATOM   2614 C   C   . THR THR A A 182 182 . -9.319  23.770 1.839   1.00 31.83 .
ATOM   2615 O   O   . THR THR A A 182 182 . -8.713  23.467 0.811   1.00 32.83 .
ATOM   2616 CB  CB  . THR THR A A 182 182 . -9.279  26.232 2.237   1.00 30.72 .
ATOM   2617 OG1 OG1 . THR THR A A 182 182 . -8.703  26.490 0.952   1.00 27.58 .
ATOM   2618 CG2 CG2 . THR THR A A 182 182 . -8.864  27.338 3.202   1.00 25.39 .
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ATOM   2620 CA  CA  . ASP ASP A A 183 183 . -11.033 22.113 1.399   1.00 30.68 .
ATOM   2621 C   C   . ASP ASP A A 183 183 . -11.448 22.693 0.054   1.00 25.46 .
ATOM   2622 O   O   . ASP ASP A A 183 183 . -11.349 22.030 -0.970  1.00 37.77 .
ATOM   2623 CB  CB  . ASP ASP A A 183 183 . -12.252 21.491 2.089   1.00 37.83 .
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ATOM   2626 OD2 OD2 . ASP ASP A A 183 183 . -12.510 20.836 4.365   1.00 57.30 .
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ATOM   2630 O   O   . SER SER A A 184 184 . -10.810 26.365 -0.679  1.00 31.39 .
ATOM   2631 CB  CB  . SER SER A A 184 184 . -13.804 24.914 -1.131  1.00 37.93 .
ATOM   2632 OG  OG  . SER SER A A 184 184 . -14.588 23.752 -0.913  1.00 48.99 .
ATOM   2633 N   N   . ILE ILE A A 185 185 . -11.603 26.248 -2.781  1.00 30.62 .
ATOM   2634 CA  CA  . ILE ILE A A 185 185 . -10.900 27.429 -3.240  1.00 34.03 .
ATOM   2635 C   C   . ILE ILE A A 185 185 . -11.574 28.663 -2.653  1.00 38.86 .
ATOM   2636 O   O   . ILE ILE A A 185 185 . -12.793 28.805 -2.723  1.00 38.25 .
ATOM   2637 CB  CB  . ILE ILE A A 185 185 . -10.927 27.512 -4.777  1.00 29.91 .
ATOM   2638 CG1 CG1 . ILE ILE A A 185 185 . -10.170 26.316 -5.367  1.00 43.46 .
ATOM   2639 CG2 CG2 . ILE ILE A A 185 185 . -10.302 28.814 -5.243  1.00 23.99 .
ATOM   2640 CD1 CD1 . ILE ILE A A 185 185 . -10.167 26.262 -6.883  1.00 40.21 .
ATOM   2641 N   N   . THR THR A A 186 186 . -10.781 29.553 -2.069  1.00 35.03 .
ATOM   2642 CA  CA  . THR THR A A 186 186 . -11.325 30.764 -1.476  1.00 33.19 .
ATOM   2643 C   C   . THR THR A A 186 186 . -10.685 32.014 -2.068  1.00 33.13 .
ATOM   2644 O   O   . THR THR A A 186 186 . -9.902  31.938 -3.023  1.00 31.78 .
ATOM   2645 CB  CB  . THR THR A A 186 186 . -11.122 30.774 0.058   1.00 34.11 .
ATOM   2646 OG1 OG1 . THR THR A A 186 186 . -9.721  30.828 0.364   1.00 30.91 .
ATOM   2647 CG2 CG2 . THR THR A A 186 186 . -11.717 29.521 0.679   1.00 30.13 .
ATOM   2648 N   N   . GLY GLY A A 187 187 . -11.033 33.164 -1.499  1.00 26.68 .
ATOM   2649 CA  CA  . GLY GLY A A 187 187 . -10.484 34.422 -1.967  1.00 25.97 .
ATOM   2650 C   C   . GLY GLY A A 187 187 . -9.044  34.592 -1.515  1.00 21.91 .
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ATOM   2663 CA  CA  . ALA ALA A A 190 190 . -4.770  32.748 -5.540  1.00 17.20 .
ATOM   2664 C   C   . ALA ALA A A 190 190 . -3.391  33.315 -5.202  1.00 19.32 .
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ATOM   2666 CB  CB  . ALA ALA A A 190 190 . -5.568  33.764 -6.331  1.00 23.09 .
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ATOM   2670 O   O   . TRP TRP A A 191 191 . -0.802  33.336 -8.026  1.00 21.60 .
ATOM   2671 CB  CB  . TRP TRP A A 191 191 . -0.127  32.136 -5.012  1.00 15.87 .
ATOM   2672 CG  CG  . TRP TRP A A 191 191 . -0.640  31.640 -3.688  1.00 13.53 .
ATOM   2673 CD1 CD1 . TRP TRP A A 191 191 . -1.576  30.667 -3.483  1.00 13.97 .
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ATOM   2684 O   O   . LYS LYS A A 192 192 . 3.342   35.147 -6.414  1.00 28.55 .
ATOM   2685 CB  CB  . LYS LYS A A 192 192 . 1.083   36.467 -8.131  1.00 18.31 .
ATOM   2686 CG  CG  . LYS LYS A A 192 192 . 1.951   37.211 -9.105  1.00 33.17 .
ATOM   2687 CD  CD  . LYS LYS A A 192 192 . 1.636   38.708 -9.077  1.00 47.63 .
ATOM   2688 CE  CE  . LYS LYS A A 192 192 . 1.866   39.284 -7.682  1.00 47.45 .
ATOM   2689 NZ  NZ  . LYS LYS A A 192 192 . 1.669   40.756 -7.613  1.00 51.73 .
ATOM   2690 N   N   . VAL VAL A A 193 193 . 3.778   34.543 -8.535  1.00 23.62 .
ATOM   2691 CA  CA  . VAL VAL A A 193 193 . 5.211   34.399 -8.319  1.00 31.91 .
ATOM   2692 C   C   . VAL VAL A A 193 193 . 5.955   35.686 -8.651  1.00 37.20 .
ATOM   2693 O   O   . VAL VAL A A 193 193 . 5.493   36.490 -9.455  1.00 41.85 .
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ATOM   2695 CG1 CG1 . VAL VAL A A 193 193 . 7.284   33.363 -9.304  1.00 44.66 .
ATOM   2696 CG2 CG2 . VAL VAL A A 193 193 . 5.429   31.908 -8.553  1.00 40.07 .
ATOM   2697 N   N   . ALA ALA A A 194 194 . 7.105   35.881 -8.016  1.00 43.72 .
ATOM   2698 CA  CA  . ALA ALA A A 194 194 . 7.932   37.059 -8.255  1.00 42.95 .
ATOM   2699 C   C   . ALA ALA A A 194 194 . 9.384   36.697 -7.962  1.00 48.32 .
ATOM   2700 O   O   . ALA ALA A A 194 194 . 9.676   35.940 -7.030  1.00 34.61 .
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ATOM   2702 N   N   . GLU GLU A A 195 195 . 10.303  37.222 -8.761  1.00 48.23 .
ATOM   2703 CA  CA  . GLU GLU A A 195 195 . 11.703  36.921 -8.528  1.00 55.29 .
ATOM   2704 C   C   . GLU GLU A A 195 195 . 12.264  37.787 -7.417  1.00 55.48 .
ATOM   2705 O   O   . GLU GLU A A 195 195 . 12.378  39.004 -7.564  1.00 61.47 .
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ATOM   2707 CG  CG  . GLU GLU A A 195 195 . 14.022  36.973 -9.536  1.00 51.64 .
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ATOM   2712 CA  CA  . GLY GLY A A 196 196 . 13.175  37.894 -5.192  1.00 65.49 .
ATOM   2713 C   C   . GLY GLY A A 196 196 . 14.590  38.222 -5.620  1.00 68.22 .
ATOM   2714 O   O   . GLY GLY A A 196 196 . 14.943  39.386 -5.829  1.00 63.76 .
ATOM   2715 N   N   . ILE ILE A A 197 197 . 15.399  37.176 -5.762  1.00 67.90 .
ATOM   2716 CA  CA  . ILE ILE A A 197 197 . 16.781  37.307 -6.198  1.00 67.52 .
ATOM   2717 C   C   . ILE ILE A A 197 197 . 17.238  35.976 -6.789  1.00 69.46 .
ATOM   2718 O   O   . ILE ILE A A 197 197 . 18.047  35.287 -6.134  1.00 71.49 .
ATOM   2719 CB  CB  . ILE ILE A A 197 197 . 17.731  37.674 -5.044  1.00 67.81 .
ATOM   2720 CG1 CG1 . ILE ILE A A 197 197 . 17.240  38.921 -4.309  1.00 69.85 .
ATOM   2721 CG2 CG2 . ILE ILE A A 197 197 . 19.103  37.956 -5.609  1.00 71.56 .
ATOM   2722 CD1 CD1 . ILE ILE A A 197 197 . 18.102  39.312 -3.123  1.00 74.69 .
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HETATM 2725 C3  C3  . MLI MLI . . 201 201 . -6.642  27.854 -12.323 1.00 44.60 .
HETATM 2726 O6  O6  . MLI MLI . . 201 201 . -8.474  25.922 -10.633 1.00 46.51 .
HETATM 2727 O7  O7  . MLI MLI . . 201 201 . -9.663  27.621 -11.085 1.00 46.46 .
HETATM 2728 O8  O8  . MLI MLI . . 201 201 . -6.222  28.726 -13.081 1.00 47.30 .
HETATM 2729 O9  O9  . MLI MLI . . 201 201 . -6.014  27.602 -11.297 1.00 56.58 .
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_atom_sites.fract_transf_matrix[1][3]   0.000000
_atom_sites.fract_transf_matrix[2][1]   0.000000
_atom_sites.fract_transf_matrix[2][2]   1.000000
_atom_sites.fract_transf_matrix[2][3]   0.000000
_atom_sites.fract_transf_matrix[3][1]   0.000000
_atom_sites.fract_transf_matrix[3][2]   0.000000
_atom_sites.fract_transf_matrix[3][3]   1.000000
_atom_sites.fract_transf_vector[1]      0.000000
_atom_sites.fract_transf_vector[2]      0.000000
_atom_sites.fract_transf_vector[3]      0.000000
#
_audit_author.name   'ProMod (SEE REFERENCE IN JRNL Records)'
#
_cell.length_a      1.000
_cell.length_b      1.000
_cell.length_c      1.000
_cell.angle_alpha   90.000
_cell.angle_beta    90.000
_cell.angle_gamma   90.000
_cell.entry_id      UNNAMED
#
_citation.id                        primary
_citation.title
_citation.journal_abbrev            (null)
_citation.journal_volume            (null)
_citation.page_first                (null)
_citation.page_last                 ?
_citation.year                      (null)
_citation.journal_id_ASTM           (null)
_citation.country                   (null)
_citation.journal_id_ISSN           (null)
_citation.journal_id_CSD            ?
_citation.book_publisher            ?
```

```
_citation.pdbx_database_id_PubMed   ?
_citation.pdbx_database_id_DOI      ?
#
_citation_author.citation_id   primary
_citation_author.name          M
_citation_author.ordinal       1
#
_computing.entry_id                     UNNAMED
_computing.structure_refinement         GROMOS96
_computing.structure_solution           ?
_computing.pdbx_density_modification    ?
_computing.pdbx_data_reduction_ds       ?
_computing.pdbx_data_reduction_ii       ?
_computing.cell_refinement              ?
#
_data_extraction.software           pdb_extract
_data_extraction.extraction_date    'Mon Aug 25 17:58:07 2008'
_data_extraction.version            3.006
_data_extraction.release_date       'June 11, 2008'
_data_extraction.location           http://sw-tools.rcsb.org/apps/PDB_EXTRACT/
#
_database_PDB_matrix.entry_id           UNNAMED
_database_PDB_matrix.origx[1][1]        1.000000
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_database_PDB_matrix.origx[1][3]        0.000000
_database_PDB_matrix.origx[2][1]        0.000000
_database_PDB_matrix.origx[2][2]        1.000000
_database_PDB_matrix.origx[2][3]        0.000000
_database_PDB_matrix.origx[3][1]        0.000000
_database_PDB_matrix.origx[3][2]        0.000000
_database_PDB_matrix.origx[3][3]        1.000000
_database_PDB_matrix.origx_vector[1]    0.000000
_database_PDB_matrix.origx_vector[2]    0.000000
_database_PDB_matrix.origx_vector[3]    0.000000
#
loop_
_database_PDB_remark.id
_database_PDB_remark.text
1
;
REMARK   1
;
2
;
REMARK   2
REMARK   2 RESOLUTION.  NOT APPLICABLE.  SEE REMARK 4.
;
3
;
REMARK   3
REMARK   3 REFINEMENT.  NONE.
REMARK   3
REMARK   3 MODELLING METHOD.
REMARK   3   PROGRAM     : ProModII
REMARK   3   AUTHORS     : N.GUEX,M.C.PEITSCH
REMARK   3
REMARK   3 ENERGY MINIMSATION.
```

```
REMARK   3   PROGRAM            : GROMOS96
REMARK   3   AUTHORS            : VAN GUNSTEREN
REMARK   3   PARAMETER          : IFP43B1
REMARK   3   TOPOLOGY FILE      : MTB43B1
REMARK   3   METHOD      1      : STEEPEST DESCENT
REMARK   3       CYCLES      1  : 200
REMARK   3       CONSTRAINTS 1  : 25 / C-FACTORS
REMARK   3   METHOD      2      : CONJUGATE GRADIENT
REMARK   3       CYCLES      2  : 300
REMARK   3       CONSTRAINTS 2  : 2500 / C-FACTORS
REMARK   3
;
4
;
REMARK   4
REMARK   4 THE COORDINATES IN THIS ENTRY REPRESENT A MODEL STRUCTURE.
REMARK   4 PROTEIN DATA BANK CONVENTIONS REQUIRE THAT *CRYST1* AND
REMARK   4 *SCALE* RECORDS BE INCLUDED, BUT THE VALUES ON THESE
REMARK   4 RECORDS ARE MEANINGLESS.
;
5
;
REMARK   5
REMARK   5 THIS MODEL IS BASED UPON THE COORDINATES OF:
REMARK   5
REMARK   5 PDB ENTRY 1t9f.pdb
REMARK   5     STRUCTURAL GENOMICS, UNKNOWN FUNCTION   16-MAY-04
REMARK   5     STRUCTURAL GENOMICS OF CAENORHABDITIS ELEGANS: STRUCTURE O
REMARK   5     A PROTEIN WITH UNKNOWN FUNCTION
REMARK   5     X-RAY DIFFRACTION
REMARK   5     RESOLUTION. 2.00 ANGSTROMS.
REMARK   5
;
#
_entity_poly.entity_id                   1
_entity_poly.pdbx_seq_one_letter_code
;KTGAELVTCGSVLKLLNTHHRVRLHSHDIKYGSGSGQQSVTGVEASDDANSYWRIRGGSE
GGCPRGSPVRCGQAVRLTHVLTGKNLHTHHFPSPLSNNQEVSAFGEDGEGDDLDLWTVRC
SGQHWEREAAVRFQHVGTSVFLSVTGEQYGSPIRGQHEVHGMPSANTHNTWKAMEGI
;
_entity_poly.pdbx_strand_id              ?
_entity_poly.type                        polypeptide(L)
_entity_poly.pdbx_target_identifier      ?
#
_exptl.crystals_number   1
_exptl.entry_id          UNNAMED
_exptl.method            'X-RAY DIFFRACTION'
#
_exptl_crystal.id                    1
_exptl_crystal.pdbx_mosaicity        ?
_exptl_crystal.pdbx_mosaicity_esd    ?
_exptl_crystal.density_Matthews      ?
_exptl_crystal.density_diffrn        ?
_exptl_crystal.density_meas          ?
_exptl_crystal.density_meas_temp     ?
_exptl_crystal.density_percent_sol   ?
_exptl_crystal.size_max              ?
```

```
_exptl_crystal.size_mid              ?
_exptl_crystal.size_min              ?
_exptl_crystal.size_rad              ?
#
_refine.entry_id                              UNNAMED
_refine.B_iso_mean                            41.006
_refine.pdbx_method_to_determine_struct       N/A
_refine.B_iso_max                             99.99
_refine.B_iso_min                             0.00
_refine.occupancy_max                         1.00
_refine.occupancy_min                         1.00
#
_reflns.entry_id   UNNAMED
#
loop_
_software.name
_software.version
_software.date
_software.type
_software.contact_author
_software.contact_author_email
_software.classification
_software.location
_software.language
_software.citation_id
REFMAC5     ?     ?               program 'Garib N. Murshudov'
garib@ysbl.york.ac.uk refinement
http://www.ccp4.ac.uk/dist/html/refmac5.html Fortran_77 1
pdb_extract 3.006 'June 11, 2008' package PDB
help@deposit.rcsb.org
'data extraction' http://sw-tools.pdb.org/apps/PDB_EXTRACT/    C++        2
#
_struct.entry_id   UNNAMED
_struct.title      SDF2-L1
#
_struct_biol.id        1
_struct_biol.details   N/A
#
_symmetry.space_group_name_H-M   'P 1'
_symmetry.entry_id               UNNAMED
#
```