# Full wwPDB X-ray Structure Validation Report (i)

Dec 8, 2023 – 03:28 am GMT

| | | |
|---|---|---|
| PDB ID | : | 1HEZ |
| Title | : | Structure of P. magnus protein L bound to a human IgM Fab. |
| Authors | : | Graille, M.; Stura, E.A.; Housden, N.G.; Bottomley, S.P.; Taussig, M.J.; Sutton, B.J.; Gore, M.G.; Charbonnier, J.B. |
| Deposited on | : | 2000-11-27 |
| Resolution | : | 2.70 Å(reported) |

The following versions of software and data (see references (i)) were used in the production of this report:

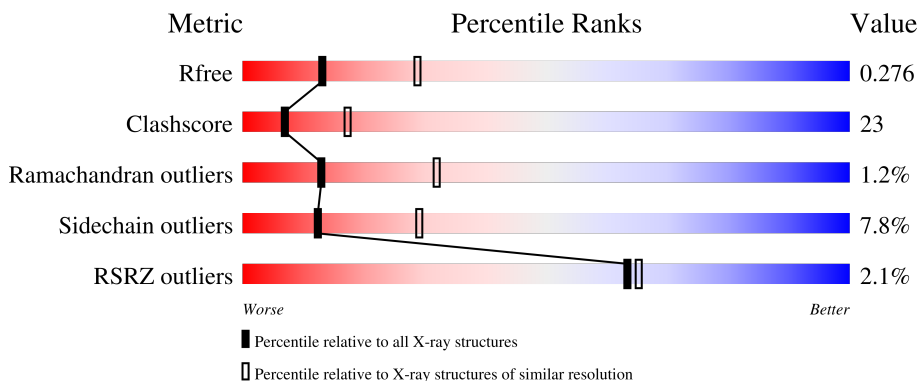| | | |
|---|---|---|
| MolProbity | : | 4.02b-467 |
| Mogul | : | 1.8.4, CSD as541be (2020) |
| Xtriage (Phenix) | : | 1.13 |
| EDS | : | 2.36 |
| Percentile statistics | : | 20191225.v01 (using entries in the PDB archive December 25th 2019) |
| Refmac | : | 5.8.0158 |
| CCP4 | : | 7.0.044 (Gargrove) |
| Ideal geometry (proteins) | : | Engh & Huber (2001) |
| Ideal geometry (DNA, RNA) | : | Parkinson et al. (1996) |
| Validation Pipeline (wwPDB-VP) | : | 2.36 |

# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.

| Metric | Percentile Ranks | Value |
|---|---|---|
| Rfree | | 0.276 |
| Clashscore | | 23 |
| Ramachandran outliers | | 1.2% |
| Sidechain outliers | | 7.8% |
| RSRZ outliers | | 2.1% |

*Worse*        *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

| Metric | Whole archive (#Entries) | Similar resolution (#Entries, resolution range(Å)) |
|---|---|---|
| R$_{free}$ | 130704 | 2808 (2.70-2.70) |
| Clashscore | 141614 | 3122 (2.70-2.70) |
| Ramachandran outliers | 138981 | 3069 (2.70-2.70) |
| Sidechain outliers | 138945 | 3069 (2.70-2.70) |
| RSRZ outliers | 127900 | 2737 (2.70-2.70) |

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

| Mol | Chain | Length | Quality of chain |
|---|---|---|---|
| 1 | A | 214 | 2%<br>56%   40%   • |
| 1 | C | 214 | 66%   30%   • |
| 2 | B | 224 | 3%<br>48%   41%   •   8% |
| 2 | D | 224 | 3%<br>55%   38%   •   5% |
| 3 | E | 61 | 2%<br>61%   31%   8% |

WORLDWIDE PDB
PROTEIN DATA BANK

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

| Mol | Type | Chain | Res | Chirality | Geometry | Clashes | Electron density |
|-----|------|-------|-----|-----------|----------|---------|------------------|
| 4 | IMD | A | 401 | - | - | X | - |
| 4 | IMD | C | 401 | - | - | X | - |

## 2  Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 7058 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called KAPPA LIGHT CHAIN OF IG.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|---|---|---|---|---------|---------|-------|
| 1 | A | 214 | Total<br>1634 | C<br>1016 | N<br>274 | O<br>338 | S<br>6 | 0 | 0 | 0 |
| 1 | C | 214 | Total<br>1635 | C<br>1016 | N<br>275 | O<br>338 | S<br>6 | 0 | 0 | 0 |

- Molecule 2 is a protein called HEAVY CHAIN OF IG.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|---|---|---|---|---------|---------|-------|
| 2 | B | 207 | Total<br>1583 | C<br>1005 | N<br>271 | O<br>300 | S<br>7 | 0 | 0 | 0 |
| 2 | D | 213 | Total<br>1616 | C<br>1023 | N<br>275 | O<br>311 | S<br>7 | 0 | 0 | 0 |

- Molecule 3 is a protein called PROTEIN L.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|---|---|---|---|---------|---------|-------|
| 3 | E | 61 | Total<br>468 | C<br>296 | N<br>74 | O<br>97 | S<br>1 | 0 | 0 | 0 |

There are 4 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|---------|-----------|
| E | 834 | ILE | THR | conflict | UNP Q51918 |
| E | 873 | ASN | TYR | conflict | UNP Q51918 |
| E | 874 | HIS | THR | conflict | UNP Q51918 |
| E | 875 | MET | ILE | conflict | UNP Q51918 |

- Molecule 4 is IMIDAZOLE (three-letter code: IMD) (formula: $C_3H_5N_2$).

IMD

N3

H

N+

C4

C2

N

C5

H

N1

| Mol | Chain | Residues | Atoms | | | ZeroOcc | AltConf |
|---|---|---|---|---|---|---|---|
| 4 | A | 1 | Total C N | | | 0 | 0 |
| | | | 5 | 3 | 2 | | |
| 4 | C | 1 | Total C N | | | 0 | 0 |
| | | | 5 | 3 | 2 | | |

- Molecule 5 is water.

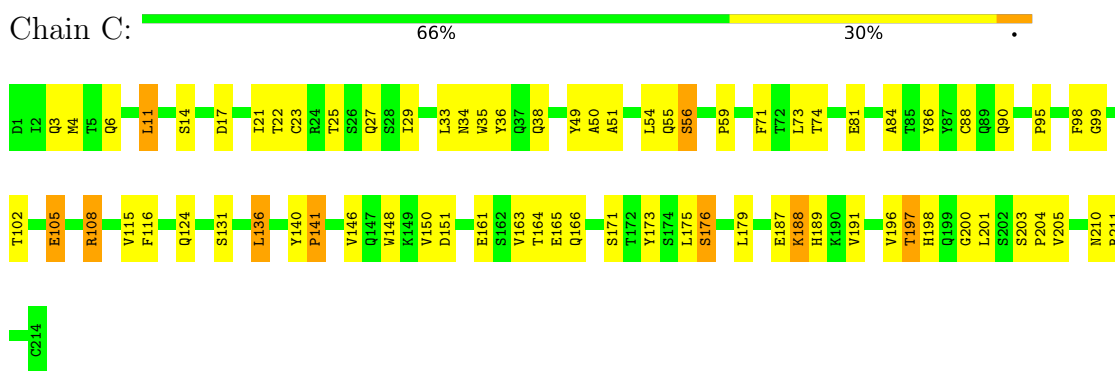| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|---|---|---|---|---|---|---|
| 5 | A | 33 | Total O | | 0 | 0 |
| | | | 33 | 33 | | |
| 5 | B | 18 | Total O | | 0 | 0 |
| | | | 18 | 18 | | |
| 5 | C | 28 | Total O | | 0 | 0 |
| | | | 28 | 28 | | |
| 5 | D | 26 | Total O | | 0 | 0 |
| | | | 26 | 26 | | |
| 5 | E | 7 | Total O | | 0 | 0 |
| | | | 7 | 7 | | |

# 3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.
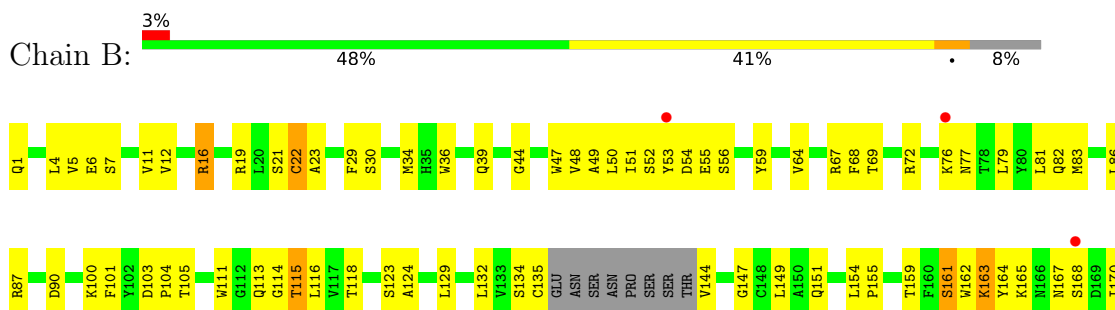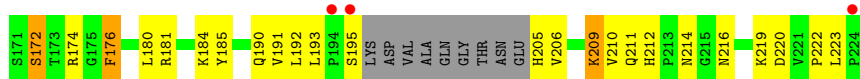
- Molecule 1: KAPPA LIGHT CHAIN OF IG

Chain A:
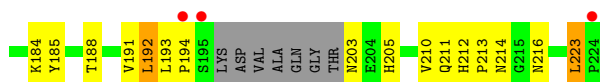


- Molecule 1: KAPPA LIGHT CHAIN OF IG

Chain C:



- Molecule 2: HEAVY CHAIN OF IG

Chain B:

S171 S172 T173 R174 G175 F176 L180 R181 K184 Y185 Q190 V191 L192 L193 P194 S195 LYS ASP VAL ALA GLN GLY THR ASN GLU H205 V206 K209 V210 Q211 H212 P213 N214 G215 N216 K219 D220 V221 P222 L223 P224

• Molecule 2: HEAVY CHAIN OF IG

Chain D:

3%

55%　　　　　　38%　　　　• 5%

Q1 V2 E6 S7 V12 Q13 P14 G15 R16 S17 L18 R19 C22 F29 M34 H35 W36 Q39 W47 L50 D54 E55 Y59 Y60 R67 R72 D73 N74 S75 K76 N77 T78 L79 Y80 L81 Q82 M83 N84 S85 L86 T91 A92 C96 A97 K100

F101 Y102 D103 P107 Y110 W111 T115 L116 V119 S120 S121 A124 S125 A126 P127 E136 N137 SER ASN PRO SER S142 T143 V144 A145 V146 G147 C148 L149 D152 F153 L154 P155 D156 S157 I158 T159 W162 K163 Y164 K165 N166 N167 I170 S171 S172 T173 F176 L180

K184 Y185 T188 V191 L192 L193 P194 S195 LYS ASP VAL ALA GLN GLY THR N203 E204 H205 V210 Q211 H212 P213 N214 G215 N216 L223 P224

• Molecule 3: PROTEIN L

Chain E:

2%

61%　　　　　　31%　　　　8%

E820 V821 T822 V825 F829 A830 D831 I834 Q835 T836 T842 E845 R852 L856 L857 A858 K859 V860 N861 G862 E863 Y864 T865 A866 D867 L868 G872 M875 N876 I877 A880

# 4 Data and refinement statistics ⓘ

| Property | Value | Source |
|---|---|---|
| Space group | P 21 21 21 | Depositor |
| Cell constants<br>a, b, c, $\alpha$, $\beta$, $\gamma$ | 55.19Å   87.33Å   210.54Å<br>90.00°   90.00°   90.00° | Depositor |
| Resolution (Å) | 20.00  –  2.70<br>19.94  –  2.60 | Depositor<br>EDS |
| % Data completeness<br>(in resolution range) | 84.5 (20.00-2.70)<br>88.3 (19.94-2.60) | Depositor<br>EDS |
| $R_{merge}$ | (Not available) | Depositor |
| $R_{sym}$ | 0.11 | Depositor |
| $< I/\sigma(I) >$ [1] | 1.13 (at 2.59Å) | Xtriage |
| Refinement program | CNS 1.0 | Depositor |
| R, $R_{free}$ | 0.215  ,  0.278<br>0.218  ,  0.276 | Depositor<br>DCC |
| $R_{free}$ test set | 1473 reflections (4.92%) | wwPDB-VP |
| Wilson B-factor (Å$^2$) | 46.1 | Xtriage |
| Anisotropy | 0.315 | Xtriage |
| Bulk solvent $k_{sol}$(e/Å$^3$), $B_{sol}$(Å$^2$) | 0.31 , 43.1 | EDS |
| L-test for twinning[2] | $< |L| > = 0.48, < L^2 > = 0.32$ | Xtriage |
| Estimated twinning fraction | No twinning to report. | Xtriage |
| $F_o$,$F_c$ correlation | 0.92 | EDS |
| Total number of atoms | 7058 | wwPDB-VP |
| Average B, all atoms (Å$^2$) | 45.0 | wwPDB-VP |

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.97% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1] Intensities estimated from amplitudes.
[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

# 5　Model quality ⓘ

## 5.1　Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: IMD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mol | Chain | Bond lengths | | Bond angles | |
|-----|-------|------|---------|------|---------|
| | | RMSZ | #$|Z|$>5 | RMSZ | #$|Z|$>5 |
| 1 | A | 0.40 | 0/1668 | 0.66 | 0/2265 |
| 1 | C | 0.42 | 0/1669 | 0.69 | 0/2265 |
| 2 | B | 0.39 | 0/1622 | 0.67 | 0/2202 |
| 2 | D | 0.42 | 0/1655 | 0.68 | 0/2249 |
| 3 | E | 0.48 | 0/475 | 0.63 | 0/641 |
| All | All | 0.41 | 0/7089 | 0.67 | 0/9622 |

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

## 5.2　Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 1 | A | 1634 | 0 | 1573 | 80 | 0 |
| 1 | C | 1635 | 0 | 1575 | 56 | 0 |
| 2 | B | 1583 | 0 | 1543 | 89 | 0 |
| 2 | D | 1616 | 0 | 1558 | 75 | 0 |
| 3 | E | 468 | 0 | 434 | 26 | 0 |
| 4 | A | 5 | 0 | 5 | 5 | 0 |
| 4 | C | 5 | 0 | 5 | 5 | 0 |
| 5 | A | 33 | 0 | 0 | 1 | 0 |

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| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 5 | B | 18 | 0 | 0 | 0 | 0 |
| 5 | C | 28 | 0 | 0 | 1 | 0 |
| 5 | D | 26 | 0 | 0 | 0 | 0 |
| 5 | E | 7 | 0 | 0 | 0 | 0 |
| All | All | 7058 | 0 | 6693 | 314 | 0 |

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 23.

All (314) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|--------|--------|--------------------------|-------------------|
| 2:D:149:LEU:HD12 | 2:D:188:THR:HG22 | 1.43 | 1.01 |
| 1:A:211:ARG:HB3 | 1:A:211:ARG:HH11 | 1.23 | 0.99 |
| 1:A:29:ILE:HG23 | 1:A:92:TYR:HB2 | 1.52 | 0.92 |
| 2:B:154:LEU:HD12 | 2:B:155:PRO:HA | 1.52 | 0.91 |
| 1:A:2:ILE:HG21 | 1:A:29:ILE:HD11 | 1.56 | 0.88 |
| 3:E:852:ARG:HH11 | 3:E:852:ARG:HB3 | 1.39 | 0.86 |
| 2:B:19:ARG:HB3 | 2:B:19:ARG:NH1 | 1.92 | 0.85 |
| 1:C:188:LYS:HA | 1:C:188:LYS:NZ | 1.92 | 0.85 |
| 1:A:98:PHE:HB2 | 4:A:401:IMD:HN3 | 1.42 | 0.84 |
| 2:D:1:GLN:HA | 2:D:1:GLN:OE1 | 1.77 | 0.84 |
| 1:A:107:LYS:HE3 | 3:E:834:ILE:HG13 | 1.62 | 0.81 |
| 1:C:6:GLN:HE21 | 1:C:99:GLY:HA3 | 1.47 | 0.80 |
| 3:E:834:ILE:HG13 | 3:E:834:ILE:O | 1.82 | 0.78 |
| 1:C:175:LEU:HD23 | 1:C:176:SER:N | 1.99 | 0.78 |
| 2:D:159:THR:HB | 2:D:211:GLN:HB2 | 1.64 | 0.78 |
| 2:B:12:VAL:HG11 | 2:B:86:LEU:HD13 | 1.66 | 0.77 |
| 3:E:852:ARG:HH11 | 3:E:852:ARG:CB | 1.97 | 0.77 |
| 2:B:181:ARG:HH11 | 2:B:181:ARG:HB3 | 1.49 | 0.76 |
| 3:E:852:ARG:HB3 | 3:E:852:ARG:NH1 | 2.00 | 0.76 |
| 2:D:158:ILE:HD11 | 2:D:210:VAL:HB | 1.67 | 0.75 |
| 2:B:206:VAL:HG13 | 2:B:223:LEU:HD21 | 1.69 | 0.74 |
| 1:C:98:PHE:H | 4:C:401:IMD:HN3 | 1.33 | 0.73 |
| 2:D:124:ALA:HB1 | 2:D:214:ASN:HD21 | 1.53 | 0.73 |
| 2:B:67:ARG:HD2 | 2:B:87:ARG:NH2 | 2.03 | 0.73 |
| 2:B:165:LYS:HE2 | 2:B:205:HIS:CE1 | 2.24 | 0.72 |
| 1:A:211:ARG:HH11 | 1:A:211:ARG:CB | 1.99 | 0.72 |
| 3:E:829:PHE:HE1 | 3:E:835:GLN:HB2 | 1.55 | 0.72 |
| 1:C:11:LEU:HD11 | 3:E:868:LEU:HD13 | 1.74 | 0.70 |
| 1:A:3:GLN:HB2 | 1:A:26:SER:HB3 | 1.73 | 0.69 |

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| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|---|---|---|---|
| 1:C:33:LEU:HD22 | 1:C:71:PHE:CG | 2.27 | 0.69 |
| 2:D:148:CYS:HB2 | 2:D:162:TRP:CH2 | 2.28 | 0.69 |
| 1:A:210:ASN:O | 1:A:213:GLU:HB2 | 1.93 | 0.69 |
| 1:A:143:GLU:HB2 | 5:A:2028:HOH:O | 1.93 | 0.69 |
| 1:A:108:ARG:HG3 | 1:A:109:THR:N | 2.07 | 0.69 |
| 4:A:401:IMD:H2 | 2:B:47:TRP:H | 1.58 | 0.68 |
| 2:D:192:LEU:N | 2:D:192:LEU:HD22 | 2.08 | 0.68 |
| 2:D:29:PHE:CD2 | 2:D:77:ASN:HA | 2.29 | 0.68 |
| 1:A:141:PRO:O | 1:A:198:HIS:HE1 | 1.77 | 0.67 |
| 1:A:29:ILE:O | 1:A:29:ILE:HG22 | 1.95 | 0.66 |
| 1:C:124:GLN:HE22 | 1:C:131:SER:N | 1.93 | 0.66 |
| 1:C:198:HIS:CD2 | 1:C:200:GLY:H | 2.14 | 0.66 |
| 2:D:143:THR:HB | 2:D:193:LEU:O | 1.97 | 0.65 |
| 1:C:73:LEU:HD23 | 1:C:74:THR:N | 2.12 | 0.65 |
| 1:A:131:SER:HB3 | 1:A:180:THR:HG23 | 1.78 | 0.65 |
| 2:B:11:VAL:HG21 | 2:B:124:ALA:HB2 | 1.79 | 0.65 |
| 1:A:29:ILE:HG23 | 1:A:92:TYR:CB | 2.27 | 0.64 |
| 1:A:198:HIS:CD2 | 1:A:200:GLY:H | 2.15 | 0.63 |
| 1:C:197:THR:HG23 | 1:C:204:PRO:HG3 | 1.81 | 0.63 |
| 1:A:136:LEU:HG | 1:A:175:LEU:HD22 | 1.81 | 0.63 |
| 1:A:158:ASN:HD22 | 1:A:181:LEU:HD21 | 1.63 | 0.63 |
| 2:B:19:ARG:HB3 | 2:B:19:ARG:HH11 | 1.63 | 0.63 |
| 2:B:118:THR:HG21 | 2:B:154:LEU:HD11 | 1.79 | 0.62 |
| 2:D:172:SER:OG | 2:D:191:VAL:HG13 | 2.00 | 0.62 |
| 3:E:825:VAL:HG22 | 3:E:875:MET:HG2 | 1.81 | 0.62 |
| 2:B:180:LEU:HD12 | 2:B:184:LYS:O | 2.00 | 0.61 |
| 2:B:1:GLN:HA | 2:B:1:GLN:OE1 | 2.00 | 0.61 |
| 1:C:136:LEU:N | 1:C:136:LEU:HD23 | 2.15 | 0.61 |
| 3:E:821:VAL:HG22 | 3:E:822:THR:N | 2.15 | 0.61 |
| 1:C:188:LYS:HA | 1:C:188:LYS:HZ3 | 1.66 | 0.60 |
| 3:E:829:PHE:HD2 | 3:E:861:ASN:OD1 | 1.84 | 0.60 |
| 2:B:163:LYS:HG3 | 2:B:168:SER:O | 2.01 | 0.60 |
| 1:A:33:LEU:HD12 | 1:A:89:GLN:O | 2.01 | 0.60 |
| 3:E:829:PHE:CE1 | 3:E:835:GLN:HB2 | 2.36 | 0.60 |
| 1:A:98:PHE:HB2 | 4:A:401:IMD:N3 | 2.16 | 0.60 |
| 1:C:151:ASP:OD2 | 1:C:189:HIS:HB3 | 2.02 | 0.59 |
| 2:D:170:ILE:HG21 | 2:D:191:VAL:HG11 | 1.84 | 0.59 |
| 2:B:129:LEU:HD21 | 2:B:210:VAL:HG21 | 1.85 | 0.59 |
| 1:C:188:LYS:HA | 1:C:188:LYS:HZ2 | 1.66 | 0.59 |
| 2:B:161:SER:HB3 | 2:B:209:LYS:NZ | 2.17 | 0.59 |
| 2:D:146:VAL:HG11 | 2:D:223:LEU:HD23 | 1.85 | 0.59 |

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| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|---|---|---|---|
| 2:B:181:ARG:HB3 | 2:B:181:ARG:NH1 | 2.16 | 0.58 |
| 2:B:50:LEU:C | 2:B:50:LEU:HD12 | 2.24 | 0.58 |
| 2:D:34:MET:HB3 | 2:D:79:LEU:HD22 | 1.86 | 0.58 |
| 1:C:140:TYR:HA | 1:C:141:PRO:O | 2.04 | 0.57 |
| 2:B:212:HIS:CD2 | 2:B:214:ASN:H | 2.22 | 0.57 |
| 1:A:175:LEU:C | 1:A:175:LEU:HD23 | 2.24 | 0.57 |
| 4:C:401:IMD:C2 | 2:D:47:TRP:H | 2.17 | 0.57 |
| 1:C:191:VAL:HG22 | 1:C:210:ASN:HD21 | 1.70 | 0.57 |
| 1:A:115:VAL:HG21 | 1:A:196:VAL:HG21 | 1.87 | 0.57 |
| 2:B:30:SER:O | 2:B:53:TYR:HB2 | 2.04 | 0.57 |
| 1:C:191:VAL:HG22 | 1:C:210:ASN:ND2 | 2.20 | 0.57 |
| 1:A:180:THR:O | 1:A:181:LEU:HD23 | 2.04 | 0.56 |
| 2:D:19:ARG:HB2 | 2:D:82:GLN:OE1 | 2.04 | 0.56 |
| 1:A:6:GLN:HG3 | 1:A:88:CYS:SG | 2.44 | 0.56 |
| 1:A:137:ASN:HB3 | 1:A:138:ASN:OD1 | 2.06 | 0.56 |
| 2:B:219:LYS:HG3 | 2:B:220:ASP:H | 1.69 | 0.56 |
| 2:D:142:SER:O | 2:D:143:THR:HG23 | 2.05 | 0.56 |
| 2:D:165:LYS:HE2 | 2:D:205:HIS:CE1 | 2.41 | 0.56 |
| 2:B:49:ALA:HB1 | 2:B:59:TYR:O | 2.05 | 0.56 |
| 2:D:22:CYS:HB3 | 2:D:79:LEU:HB3 | 1.88 | 0.56 |
| 2:D:159:THR:O | 2:D:210:VAL:HA | 2.06 | 0.56 |
| 1:A:33:LEU:HD22 | 1:A:71:PHE:CD1 | 2.41 | 0.56 |
| 1:A:211:ARG:HB3 | 1:A:211:ARG:NH1 | 2.08 | 0.56 |
| 1:A:37:GLN:HB2 | 1:A:47:LEU:HD11 | 1.87 | 0.55 |
| 2:B:16:ARG:HG3 | 2:B:16:ARG:HH11 | 1.71 | 0.55 |
| 2:D:211:GLN:HG2 | 2:D:216:ASN:ND2 | 2.21 | 0.55 |
| 2:B:209:LYS:HZ3 | 2:B:211:GLN:HG3 | 1.72 | 0.55 |
| 1:C:6:GLN:HE21 | 1:C:99:GLY:CA | 2.17 | 0.55 |
| 2:D:180:LEU:HD12 | 2:D:184:LYS:O | 2.06 | 0.55 |
| 2:B:19:ARG:HB3 | 2:B:19:ARG:CZ | 2.37 | 0.55 |
| 1:A:105:GLU:OE2 | 1:A:173:TYR:OH | 2.25 | 0.54 |
| 1:A:29:ILE:HD13 | 1:A:90:GLN:HB2 | 1.90 | 0.54 |
| 1:C:6:GLN:HB3 | 5:C:2002:HOH:O | 2.07 | 0.54 |
| 2:D:164:TYR:HB2 | 2:D:166:ASN:ND2 | 2.23 | 0.54 |
| 2:D:100:LYS:HD3 | 2:D:101:PHE:CZ | 2.43 | 0.54 |
| 2:D:67:ARG:HB3 | 2:D:84:ASN:O | 2.08 | 0.54 |
| 1:A:25:THR:OG1 | 1:A:69:THR:HA | 2.08 | 0.54 |
| 1:A:108:ARG:HG3 | 1:A:109:THR:H | 1.70 | 0.54 |
| 1:A:100:GLN:OE1 | 1:A:100:GLN:N | 2.38 | 0.54 |
| 4:A:401:IMD:C2 | 2:B:47:TRP:H | 2.21 | 0.54 |
| 2:B:53:TYR:HA | 2:B:72:ARG:NH1 | 2.23 | 0.54 |

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| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|---|---|---|---|
| 2:B:103:ASP:OD1 | 2:B:104:PRO:HD2 | 2.08 | 0.53 |
| 2:B:6:GLU:OE2 | 2:B:114:GLY:N | 2.39 | 0.53 |
| 1:A:125:LEU:CD1 | 1:A:130:ALA:HB2 | 2.39 | 0.53 |
| 2:B:161:SER:HB3 | 2:B:209:LYS:CE | 2.39 | 0.53 |
| 3:E:857:LEU:O | 3:E:861:ASN:HB2 | 2.09 | 0.53 |
| 2:B:162:TRP:HB3 | 2:B:170:ILE:HD12 | 1.91 | 0.53 |
| 1:A:131:SER:HB2 | 1:A:178:THR:CG2 | 2.39 | 0.52 |
| 2:B:149:LEU:HG | 2:B:151:GLN:HG2 | 1.90 | 0.52 |
| 1:A:38:GLN:O | 1:A:84:ALA:HB1 | 2.10 | 0.52 |
| 1:C:163:VAL:HG22 | 1:C:164:THR:O | 2.10 | 0.52 |
| 2:D:158:ILE:HD11 | 2:D:210:VAL:CB | 2.39 | 0.52 |
| 1:A:25:THR:HG21 | 1:A:29:ILE:HD12 | 1.92 | 0.52 |
| 2:B:159:THR:O | 2:B:210:VAL:HA | 2.10 | 0.52 |
| 2:B:172:SER:OG | 2:B:192:LEU:HB2 | 2.10 | 0.52 |
| 2:D:1:GLN:OE1 | 2:D:1:GLN:CA | 2.54 | 0.52 |
| 2:D:14:PRO:HG3 | 2:D:119:VAL:HG12 | 1.91 | 0.52 |
| 1:A:142:ARG:HB2 | 1:A:173:TYR:CE2 | 2.45 | 0.52 |
| 2:B:181:ARG:HH11 | 2:B:181:ARG:CB | 2.21 | 0.52 |
| 2:D:124:ALA:HB1 | 2:D:214:ASN:ND2 | 2.22 | 0.52 |
| 2:D:55:GLU:HG2 | 2:D:72:ARG:HG2 | 1.92 | 0.52 |
| 1:C:124:GLN:HE22 | 1:C:131:SER:H | 1.55 | 0.52 |
| 3:E:863:GLU:HA | 3:E:863:GLU:OE2 | 2.09 | 0.51 |
| 1:A:198:HIS:HD2 | 1:A:200:GLY:H | 1.56 | 0.51 |
| 2:B:172:SER:O | 2:B:191:VAL:HA | 2.10 | 0.51 |
| 2:D:36:TRP:NE1 | 2:D:81:LEU:HB2 | 2.24 | 0.51 |
| 1:A:90:GLN:HE21 | 1:A:97:THR:H | 1.56 | 0.51 |
| 1:A:108:ARG:CG | 1:A:109:THR:N | 2.72 | 0.51 |
| 2:B:67:ARG:NH2 | 2:B:90:ASP:OD2 | 2.42 | 0.51 |
| 1:C:115:VAL:HG12 | 1:C:116:PHE:N | 2.25 | 0.51 |
| 1:A:11:LEU:HD12 | 3:E:836:THR:O | 2.11 | 0.51 |
| 3:E:825:VAL:HG22 | 3:E:875:MET:CG | 2.41 | 0.51 |
| 1:A:4:MET:SD | 1:A:25:THR:HG22 | 2.51 | 0.51 |
| 2:D:154:LEU:HD12 | 2:D:155:PRO:HA | 1.92 | 0.51 |
| 2:B:64:VAL:HB | 2:B:68:PHE:CG | 2.45 | 0.50 |
| 2:D:13:GLN:HE21 | 2:D:121:SER:HA | 1.76 | 0.50 |
| 1:A:108:ARG:NH1 | 1:A:109:THR:O | 2.44 | 0.50 |
| 2:D:127:PRO:HB3 | 2:D:153:PHE:CD2 | 2.46 | 0.50 |
| 3:E:842:THR:OG1 | 3:E:845:GLU:HG3 | 2.12 | 0.50 |
| 1:A:188:LYS:HE3 | 1:A:188:LYS:HA | 1.94 | 0.50 |
| 2:B:219:LYS:HG3 | 2:B:220:ASP:N | 2.26 | 0.50 |
| 2:D:35:HIS:CD2 | 2:D:50:LEU:HB3 | 2.47 | 0.50 |

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| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|---|---|---|---|
| 2:B:6:GLU:HA | 2:B:21:SER:O | 2.11 | 0.50 |
| 2:B:176:PHE:CD1 | 2:B:176:PHE:N | 2.79 | 0.50 |
| 2:B:190:GLN:HE21 | 2:B:192:LEU:HD21 | 1.76 | 0.50 |
| 2:D:35:HIS:CD2 | 2:D:50:LEU:CB | 2.94 | 0.50 |
| 1:A:32:TYR:CD2 | 2:B:105:THR:HB | 2.46 | 0.50 |
| 2:B:180:LEU:HD13 | 2:B:185:TYR:CE1 | 2.47 | 0.50 |
| 1:A:148:TRP:CE3 | 1:A:179:LEU:HD22 | 2.48 | 0.49 |
| 2:D:165:LYS:C | 2:D:167:ASN:H | 2.14 | 0.49 |
| 2:B:144:VAL:HG22 | 2:B:193:LEU:O | 2.13 | 0.49 |
| 1:C:201:LEU:HD13 | 1:C:205:VAL:HG23 | 1.94 | 0.49 |
| 1:A:43:ALA:H | 2:B:113:GLN:HE22 | 1.59 | 0.49 |
| 2:B:118:THR:HG21 | 2:B:154:LEU:CD1 | 2.43 | 0.49 |
| 1:C:22:THR:CG2 | 1:C:23:CYS:N | 2.75 | 0.49 |
| 1:A:108:ARG:HG3 | 1:A:108:ARG:HH11 | 1.77 | 0.49 |
| 1:C:198:HIS:HD2 | 1:C:200:GLY:H | 1.61 | 0.49 |
| 1:C:21:ILE:HG21 | 1:C:102:THR:HG21 | 1.95 | 0.49 |
| 2:B:190:GLN:NE2 | 2:B:192:LEU:HD21 | 2.28 | 0.49 |
| 1:C:11:LEU:HA | 3:E:866:ALA:O | 2.13 | 0.49 |
| 1:C:98:PHE:HD1 | 4:C:401:IMD:H2 | 1.78 | 0.49 |
| 3:E:820:GLU:HG3 | 3:E:821:VAL:N | 2.27 | 0.48 |
| 1:A:129:THR:HG22 | 1:A:130:ALA:N | 2.28 | 0.48 |
| 2:B:155:PRO:O | 2:B:212:HIS:HE1 | 1.96 | 0.48 |
| 1:C:98:PHE:N | 4:C:401:IMD:HN3 | 2.08 | 0.48 |
| 1:C:49:TYR:CB | 2:D:107:PRO:HG3 | 2.43 | 0.48 |
| 2:B:67:ARG:HD2 | 2:B:87:ARG:HH21 | 1.76 | 0.48 |
| 1:A:182:SER:C | 1:A:184:ALA:N | 2.68 | 0.48 |
| 2:D:166:ASN:C | 2:D:166:ASN:HD22 | 2.16 | 0.47 |
| 1:A:98:PHE:HD1 | 4:A:401:IMD:H2 | 1.78 | 0.47 |
| 2:D:12:VAL:HG12 | 2:D:13:GLN:N | 2.30 | 0.47 |
| 1:A:29:ILE:O | 1:A:29:ILE:CG2 | 2.62 | 0.47 |
| 1:A:159:SER:HA | 1:A:178:THR:O | 2.14 | 0.47 |
| 2:B:5:VAL:O | 2:B:22:CYS:HA | 2.15 | 0.47 |
| 1:C:73:LEU:HD23 | 1:C:73:LEU:C | 2.34 | 0.47 |
| 1:C:148:TRP:CE2 | 1:C:179:LEU:HB2 | 2.49 | 0.47 |
| 1:C:55:GLN:HG3 | 1:C:56:SER:N | 2.29 | 0.47 |
| 2:D:54:ASP:O | 2:D:55:GLU:HB2 | 2.14 | 0.47 |
| 2:D:126:ALA:HB1 | 2:D:127:PRO:HD2 | 1.97 | 0.47 |
| 2:D:82:GLN:OE1 | 2:D:82:GLN:HA | 2.15 | 0.47 |
| 1:A:140:TYR:CE2 | 1:A:141:PRO:HG3 | 2.50 | 0.47 |
| 1:A:170:ASP:O | 1:A:172:THR:HG23 | 2.15 | 0.46 |
| 3:E:860:VAL:O | 3:E:860:VAL:HG23 | 2.14 | 0.46 |

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| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|---|---|---|---|
| 2:B:52:SER:O | 2:B:72:ARG:NH1 | 2.47 | 0.46 |
| 2:D:212:HIS:CD2 | 2:D:214:ASN:H | 2.33 | 0.46 |
| 1:C:33:LEU:HD22 | 1:C:71:PHE:CB | 2.45 | 0.46 |
| 2:D:111:TRP:N | 2:D:111:TRP:CD1 | 2.83 | 0.46 |
| 1:A:35:TRP:CE2 | 1:A:73:LEU:HB2 | 2.50 | 0.46 |
| 1:A:107:LYS:HE3 | 3:E:834:ILE:O | 2.16 | 0.46 |
| 2:B:50:LEU:HD12 | 2:B:51:ILE:N | 2.31 | 0.46 |
| 1:C:6:GLN:HG2 | 1:C:88:CYS:SG | 2.55 | 0.46 |
| 2:D:97:ALA:HA | 2:D:110:TYR:O | 2.16 | 0.46 |
| 2:B:223:LEU:N | 2:B:223:LEU:HD22 | 2.31 | 0.46 |
| 1:A:35:TRP:CZ3 | 1:A:88:CYS:HB3 | 2.51 | 0.46 |
| 1:A:158:ASN:ND2 | 1:A:179:LEU:HD11 | 2.30 | 0.46 |
| 1:C:176:SER:HB3 | 2:D:176:PHE:CE2 | 2.51 | 0.46 |
| 1:A:151:ASP:O | 1:A:152:ASN:HB2 | 2.15 | 0.45 |
| 2:D:148:CYS:HB2 | 2:D:162:TRP:CZ2 | 2.51 | 0.45 |
| 2:D:164:TYR:CD2 | 2:D:170:ILE:HD11 | 2.51 | 0.45 |
| 1:A:118:PHE:HE2 | 1:A:135:LEU:HD12 | 1.81 | 0.45 |
| 1:A:135:LEU:C | 1:A:136:LEU:HD23 | 2.36 | 0.45 |
| 1:C:6:GLN:NE2 | 1:C:88:CYS:H | 2.14 | 0.45 |
| 2:D:126:ALA:HB2 | 2:D:214:ASN:O | 2.17 | 0.45 |
| 2:D:124:ALA:CB | 2:D:214:ASN:HD21 | 2.27 | 0.45 |
| 2:B:34:MET:HB3 | 2:B:79:LEU:HD22 | 1.99 | 0.45 |
| 1:A:12:SER:OG | 1:A:105:GLU:HG2 | 2.17 | 0.45 |
| 1:A:62:PHE:CE2 | 1:A:75:ILE:HG12 | 2.51 | 0.45 |
| 2:B:4:LEU:HA | 2:B:23:ALA:O | 2.17 | 0.45 |
| 2:B:191:VAL:O | 2:B:191:VAL:HG12 | 2.16 | 0.45 |
| 2:B:212:HIS:HD2 | 2:B:214:ASN:H | 1.63 | 0.45 |
| 3:E:868:LEU:HG | 3:E:872:GLY:HA2 | 1.97 | 0.45 |
| 2:B:100:LYS:HD3 | 2:B:101:PHE:CE2 | 2.52 | 0.45 |
| 1:A:135:LEU:HD23 | 1:A:176:SER:HB2 | 1.99 | 0.45 |
| 2:B:11:VAL:CG2 | 2:B:124:ALA:HB2 | 2.46 | 0.45 |
| 2:B:163:LYS:HE2 | 2:B:168:SER:C | 2.37 | 0.45 |
| 2:B:205:HIS:HA | 2:B:223:LEU:CD2 | 2.47 | 0.45 |
| 1:C:150:VAL:O | 1:C:151:ASP:HB2 | 2.16 | 0.44 |
| 2:B:36:TRP:O | 2:B:48:VAL:HB | 2.17 | 0.44 |
| 2:D:83:MET:HB3 | 2:D:86:LEU:HD21 | 1.98 | 0.44 |
| 2:D:91:THR:O | 2:D:92:ALA:HB2 | 2.17 | 0.44 |
| 2:D:146:VAL:HG11 | 2:D:223:LEU:CD2 | 2.46 | 0.44 |
| 1:C:166:GLN:HG3 | 1:C:173:TYR:CZ | 2.52 | 0.44 |
| 2:B:132:LEU:HB2 | 2:B:147:GLY:C | 2.38 | 0.44 |
| 3:E:821:VAL:CG2 | 3:E:822:THR:N | 2.80 | 0.44 |

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| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|---|---|---|---|
| 1:A:161:GLU:HA | 1:A:176:SER:O | 2.17 | 0.44 |
| 1:C:146:VAL:HG23 | 1:C:146:VAL:O | 2.16 | 0.44 |
| 2:B:164:TYR:CE1 | 2:B:170:ILE:HG12 | 2.53 | 0.44 |
| 2:D:158:ILE:HA | 2:D:211:GLN:O | 2.18 | 0.44 |
| 2:B:69:THR:OG1 | 2:B:82:GLN:HB3 | 2.18 | 0.43 |
| 2:B:205:HIS:CA | 2:B:223:LEU:HD23 | 2.47 | 0.43 |
| 1:C:50:ALA:O | 1:C:51:ALA:HB3 | 2.17 | 0.43 |
| 1:C:98:PHE:HB2 | 4:C:401:IMD:HN3 | 1.82 | 0.43 |
| 1:A:137:ASN:HB3 | 2:B:174:ARG:HH12 | 1.83 | 0.43 |
| 1:C:140:TYR:CD1 | 1:C:141:PRO:HA | 2.54 | 0.43 |
| 1:A:79:GLN:HB3 | 1:A:81:GLU:OE2 | 2.18 | 0.43 |
| 2:D:73:ASP:CG | 2:D:76:LYS:HG3 | 2.38 | 0.43 |
| 1:C:95:PRO:HB3 | 2:D:47:TRP:CZ3 | 2.54 | 0.43 |
| 1:A:108:ARG:HG3 | 1:A:108:ARG:NH1 | 2.32 | 0.43 |
| 2:B:205:HIS:HA | 2:B:222:PRO:HA | 2.00 | 0.43 |
| 1:C:38:GLN:O | 1:C:84:ALA:HB1 | 2.17 | 0.43 |
| 1:A:106:ILE:H | 1:A:166:GLN:HE22 | 1.67 | 0.43 |
| 2:B:48:VAL:O | 2:B:49:ALA:HB2 | 2.19 | 0.43 |
| 1:A:209:PHE:O | 1:A:209:PHE:CD1 | 2.71 | 0.43 |
| 1:C:29:ILE:HD11 | 1:C:33:LEU:HB2 | 2.00 | 0.43 |
| 2:D:156:ASP:HB2 | 2:D:185:TYR:CG | 2.54 | 0.43 |
| 1:C:210:ASN:O | 1:C:211:ARG:C | 2.57 | 0.43 |
| 2:D:16:ARG:HG3 | 2:D:16:ARG:HH11 | 1.84 | 0.43 |
| 2:D:35:HIS:CD2 | 2:D:50:LEU:HB2 | 2.54 | 0.43 |
| 2:D:39:GLN:O | 2:D:92:ALA:HB1 | 2.18 | 0.43 |
| 3:E:875:MET:CE | 3:E:877:ILE:HD11 | 2.48 | 0.43 |
| 1:A:59:PRO:O | 1:A:61:ARG:N | 2.52 | 0.42 |
| 2:B:36:TRP:CE2 | 2:B:81:LEU:HB2 | 2.54 | 0.42 |
| 2:B:68:PHE:CE2 | 2:B:83:MET:HG2 | 2.54 | 0.42 |
| 2:B:100:LYS:HD3 | 2:B:101:PHE:HE2 | 1.83 | 0.42 |
| 2:B:211:GLN:HG2 | 2:B:216:ASN:ND2 | 2.34 | 0.42 |
| 1:C:36:TYR:O | 1:C:86:TYR:HA | 2.19 | 0.42 |
| 3:E:834:ILE:O | 3:E:834:ILE:CG1 | 2.61 | 0.42 |
| 1:C:136:LEU:HD22 | 1:C:196:VAL:HG21 | 2.01 | 0.42 |
| 2:B:129:LEU:HD11 | 2:B:210:VAL:HG22 | 2.01 | 0.42 |
| 2:D:18:LEU:HB3 | 2:D:83:MET:HE3 | 2.01 | 0.42 |
| 2:B:7:SER:HA | 2:B:115:THR:CG2 | 2.49 | 0.42 |
| 2:D:115:THR:O | 2:D:115:THR:HG23 | 2.19 | 0.42 |
| 2:D:212:HIS:HD2 | 2:D:214:ASN:H | 1.66 | 0.42 |
| 2:B:29:PHE:CD2 | 2:B:77:ASN:HA | 2.54 | 0.42 |
| 2:D:158:ILE:HG12 | 2:D:159:THR:N | 2.34 | 0.42 |

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| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|---|---|---|---|
| 2:D:165:LYS:HD3 | 2:D:203:ASN:OD1 | 2.18 | 0.42 |
| 2:B:132:LEU:HB2 | 2:B:147:GLY:CA | 2.49 | 0.42 |
| 3:E:856:LEU:O | 3:E:859:LYS:HB2 | 2.19 | 0.42 |
| 2:B:161:SER:HB3 | 2:B:209:LYS:HG3 | 2.02 | 0.42 |
| 2:D:6:GLU:O | 2:D:7:SER:HB3 | 2.19 | 0.42 |
| 1:A:186:TYR:O | 1:A:192:TYR:OH | 2.36 | 0.42 |
| 2:B:55:GLU:OE1 | 2:B:72:ARG:HG2 | 2.19 | 0.42 |
| 2:D:60:TYR:CD1 | 2:D:60:TYR:N | 2.87 | 0.42 |
| 1:A:182:SER:C | 1:A:184:ALA:H | 2.21 | 0.42 |
| 1:A:107:LYS:HE3 | 3:E:834:ILE:CG1 | 2.43 | 0.42 |
| 2:D:2:VAL:HG11 | 2:D:110:TYR:CD2 | 2.54 | 0.42 |
| 1:A:2:ILE:HG12 | 1:A:27:GLN:HG2 | 2.01 | 0.41 |
| 1:A:129:THR:CG2 | 1:A:130:ALA:N | 2.83 | 0.41 |
| 1:C:105:GLU:HB2 | 1:C:166:GLN:OE1 | 2.20 | 0.41 |
| 2:B:54:ASP:C | 2:B:56:SER:H | 2.23 | 0.41 |
| 2:B:161:SER:HB3 | 2:B:209:LYS:HE3 | 2.01 | 0.41 |
| 2:B:19:ARG:HH21 | 2:B:21:SER:HG | 1.66 | 0.41 |
| 2:D:212:HIS:CG | 2:D:213:PRO:HD2 | 2.55 | 0.41 |
| 2:B:111:TRP:N | 2:B:111:TRP:CD1 | 2.88 | 0.41 |
| 1:C:35:TRP:CE3 | 1:C:73:LEU:HD12 | 2.55 | 0.41 |
| 1:A:8:PRO:O | 1:A:102:THR:HG23 | 2.20 | 0.41 |
| 1:C:4:MET:SD | 1:C:25:THR:HG22 | 2.61 | 0.41 |
| 2:D:164:TYR:HB2 | 2:D:166:ASN:HD21 | 1.84 | 0.41 |
| 2:B:39:GLN:HG3 | 2:B:44:GLY:O | 2.20 | 0.41 |
| 2:D:35:HIS:O | 2:D:96:CYS:HA | 2.20 | 0.41 |
| 2:B:51:ILE:HD11 | 2:B:55:GLU:HB3 | 2.02 | 0.41 |
| 1:C:108:ARG:HD3 | 1:C:171:SER:HB2 | 2.02 | 0.41 |
| 1:C:146:VAL:HG22 | 1:C:161:GLU:OE1 | 2.21 | 0.41 |
| 2:D:127:PRO:HB3 | 2:D:153:PHE:HB3 | 2.03 | 0.41 |
| 1:A:182:SER:O | 1:A:185:ASP:N | 2.48 | 0.40 |
| 2:D:50:LEU:HG | 2:D:59:TYR:HB2 | 2.02 | 0.40 |
| 2:B:118:THR:CG2 | 2:B:154:LEU:HD11 | 2.50 | 0.40 |
| 2:B:134:SER:OG | 2:B:135:CYS:N | 2.54 | 0.40 |
| 1:C:33:LEU:HG | 1:C:34:ASN:N | 2.35 | 0.40 |
| 2:D:36:TRP:CE2 | 2:D:81:LEU:HB2 | 2.56 | 0.40 |
| 1:A:48:ILE:HG12 | 1:A:54:LEU:HD22 | 2.03 | 0.40 |
| 2:D:149:LEU:HD12 | 2:D:188:THR:CG2 | 2.32 | 0.40 |
| 2:B:162:TRP:HB2 | 2:B:170:ILE:HB | 2.03 | 0.40 |
| 1:C:187:GLU:C | 1:C:188:LYS:HD2 | 2.41 | 0.40 |

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed | Favoured | Allowed | Outliers | Percentiles | |
|-----|-------|----------|----------|---------|----------|----|----|
| 1 | A | 212/214 (99%) | 195 (92%) | 14 (7%) | 3 (1%) | 11 | 28 |
| 1 | C | 212/214 (99%) | 196 (92%) | 14 (7%) | 2 (1%) | 17 | 40 |
| 2 | B | 201/224 (90%) | 185 (92%) | 13 (6%) | 3 (2%) | 10 | 26 |
| 2 | D | 207/224 (92%) | 191 (92%) | 14 (7%) | 2 (1%) | 15 | 37 |
| 3 | E | 59/61 (97%) | 53 (90%) | 5 (8%) | 1 (2%) | 9 | 23 |
| All | All | 891/937 (95%) | 820 (92%) | 60 (7%) | 11 (1%) | 13 | 32 |

All (11) Ramachandran outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 2 | B | 167 | ASN |
| 1 | A | 60 | SER |
| 1 | A | 138 | ASN |
| 2 | B | 172 | SER |
| 2 | B | 16 | ARG |
| 2 | D | 136 | GLU |
| 2 | D | 194 | PRO |
| 3 | E | 831 | ASP |
| 1 | C | 59 | PRO |
| 1 | A | 84 | ALA |
| 1 | C | 141 | PRO |

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|---|---|---|---|---|---|---|
| 1 | A | 188/190 (99%) | 171 (91%) | 17 (9%) | 9 | 22 |
| 1 | C | 188/190 (99%) | 171 (91%) | 17 (9%) | 9 | 22 |
| 2 | B | 173/189 (92%) | 163 (94%) | 10 (6%) | 20 | 43 |
| 2 | D | 176/189 (93%) | 166 (94%) | 10 (6%) | 20 | 44 |
| 3 | E | 45/47 (96%) | 39 (87%) | 6 (13%) | 4 | 9 |
| All | All | 770/805 (96%) | 710 (92%) | 60 (8%) | 12 | 29 |

All (60) residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|---|---|---|---|
| 1 | A | 1 | ASP |
| 1 | A | 14 | SER |
| 1 | A | 40 | PRO |
| 1 | A | 46 | LEU |
| 1 | A | 81 | GLU |
| 1 | A | 103 | LYS |
| 1 | A | 108 | ARG |
| 1 | A | 138 | ASN |
| 1 | A | 143 | GLU |
| 1 | A | 162 | SER |
| 1 | A | 166 | GLN |
| 1 | A | 176 | SER |
| 1 | A | 180 | THR |
| 1 | A | 202 | SER |
| 1 | A | 203 | SER |
| 1 | A | 211 | ARG |
| 1 | A | 213 | GLU |
| 2 | B | 22 | CYS |
| 2 | B | 76 | LYS |
| 2 | B | 115 | THR |
| 2 | B | 116 | LEU |
| 2 | B | 123 | SER |
| 2 | B | 161 | SER |
| 2 | B | 163 | LYS |
| 2 | B | 176 | PHE |
| 2 | B | 195 | SER |
| 2 | B | 209 | LYS |
| 1 | C | 3 | GLN |
| 1 | C | 11 | LEU |
| 1 | C | 14 | SER |
| 1 | C | 17 | ASP |

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| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | C | 27 | GLN |
| 1 | C | 54 | LEU |
| 1 | C | 56 | SER |
| 1 | C | 81 | GLU |
| 1 | C | 90 | GLN |
| 1 | C | 105 | GLU |
| 1 | C | 108 | ARG |
| 1 | C | 136 | LEU |
| 1 | C | 165 | GLU |
| 1 | C | 176 | SER |
| 1 | C | 188 | LYS |
| 1 | C | 197 | THR |
| 1 | C | 203 | SER |
| 2 | D | 1 | GLN |
| 2 | D | 75 | SER |
| 2 | D | 103 | ASP |
| 2 | D | 116 | LEU |
| 2 | D | 144 | VAL |
| 2 | D | 152 | ASP |
| 2 | D | 165 | LYS |
| 2 | D | 166 | ASN |
| 2 | D | 192 | LEU |
| 2 | D | 223 | LEU |
| 3 | E | 820 | GLU |
| 3 | E | 836 | THR |
| 3 | E | 852 | ARG |
| 3 | E | 860 | VAL |
| 3 | E | 865 | THR |
| 3 | E | 868 | LEU |

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | A | 90 | GLN |
| 1 | A | 124 | GLN |
| 1 | A | 137 | ASN |
| 1 | A | 166 | GLN |
| 1 | A | 198 | HIS |
| 1 | A | 199 | GLN |
| 2 | B | 77 | ASN |
| 2 | B | 167 | ASN |
| 2 | B | 190 | GLN |

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| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 2 | B | 205 | HIS |
| 2 | B | 211 | GLN |
| 2 | B | 212 | HIS |
| 1 | C | 6 | GLN |
| 1 | C | 27 | GLN |
| 1 | C | 124 | GLN |
| 1 | C | 198 | HIS |
| 1 | C | 199 | GLN |
| 1 | C | 210 | ASN |
| 2 | D | 13 | GLN |
| 2 | D | 166 | ASN |
| 2 | D | 167 | ASN |
| 2 | D | 205 | HIS |
| 2 | D | 212 | HIS |
| 2 | D | 214 | ASN |
| 3 | E | 826 | ASN |

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mol | Type | Chain | Res | Link | Bond lengths | | | Bond angles | | |
|-----|------|-------|-----|------|--------|------|-----------|--------|------|-----------|
| | | | | | Counts | RMSZ | $\#|Z| > 2$ | Counts | RMSZ | $\#|Z| > 2$ |
| 4 | IMD | C | 401 | - | 3,5,5 | 0.60 | 0 | 4,5,5 | 0.45 | 0 |
| 4 | IMD | A | 401 | - | 3,5,5 | 0.66 | 0 | 4,5,5 | 0.48 | 0 |

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

| Mol | Type | Chain | Res | Link | Chirals | Torsions | Rings |
|-----|------|-------|-----|------|---------|----------|---------|
| 4 | IMD | C | 401 | - | - | - | 0/1/1/1 |
| 4 | IMD | A | 401 | - | - | - | 0/1/1/1 |

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 10 short contacts:

| Mol | Chain | Res | Type | Clashes | Symm-Clashes |
|-----|-------|-----|------|---------|--------------|
| 4 | C | 401 | IMD | 5 | 0 |
| 4 | A | 401 | IMD | 5 | 0 |

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

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wwPDB
PROTEIN DATA BANK

# 6 Fit of model and data ⓘ

## 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, $95^{th}$ percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

| Mol | Chain | Analysed | <RSRZ> | #RSRZ>2 | | | OWAB(Å$^2$) | Q<0.9 |
|-----|-------|----------|--------|---------|------|------|-------------|-------|
| 1 | A | 214/214 (100%) | -0.16 | 5 (2%) | 60 | 62 | 26, 43, 79, 91 | 0 |
| 1 | C | 214/214 (100%) | -0.24 | 0 | 100 | 100 | 19, 39, 56, 61 | 0 |
| 2 | B | 207/224 (92%) | 0.02 | 6 (2%) | 51 | 52 | 25, 50, 81, 92 | 0 |
| 2 | D | 213/224 (95%) | -0.16 | 7 (3%) | 46 | 46 | 20, 43, 77, 91 | 0 |
| 3 | E | 61/61 (100%) | -0.20 | 1 (1%) | 72 | 74 | 27, 40, 57, 63 | 0 |
| All | All | 909/937 (97%) | -0.14 | 19 (2%) | 63 | 65 | 19, 43, 76, 92 | 0 |

All (19) RSRZ outliers are listed below:

| Mol | Chain | Res | Type | RSRZ |
|-----|-------|-----|------|------|
| 2 | D | 195 | SER | 4.0 |
| 2 | B | 224 | PRO | 3.4 |
| 2 | B | 194 | PRO | 3.3 |
| 2 | D | 137 | ASN | 2.9 |
| 2 | D | 173 | THR | 2.8 |
| 2 | B | 195 | SER | 2.7 |
| 2 | B | 76 | LYS | 2.6 |
| 2 | B | 53 | TYR | 2.6 |
| 3 | E | 880 | ALA | 2.6 |
| 2 | B | 168 | SER | 2.5 |
| 1 | A | 127 | SER | 2.4 |
| 2 | D | 136 | GLU | 2.4 |
| 2 | D | 172 | SER | 2.4 |
| 1 | A | 184 | ALA | 2.4 |
| 2 | D | 194 | PRO | 2.3 |
| 1 | A | 214 | CYS | 2.3 |
| 1 | A | 188 | LYS | 2.2 |
| 2 | D | 224 | PRO | 2.1 |
| 1 | A | 213 | GLU | 2.0 |

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, $95^{th}$ percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

| Mol | Type | Chain | Res | Atoms | RSCC | RSR | B-factors($\text{Å}^2$) | Q<0.9 |
|-----|------|-------|-----|-------|------|-----|-------------------------|-------|
| 4 | IMD | C | 401 | 5/5 | 0.86 | 0.41 | 66,66,67,67 | 0 |
| 4 | IMD | A | 401 | 5/5 | 0.87 | 0.34 | 56,56,57,58 | 0 |

## 6.5 Other polymers ⓘ

There are no such residues in this entry.