# W O R L D W I D E
# wwPDB
## PROTEIN DATA BANK

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

Feb 27, 2023 – 06:22 pm GMT

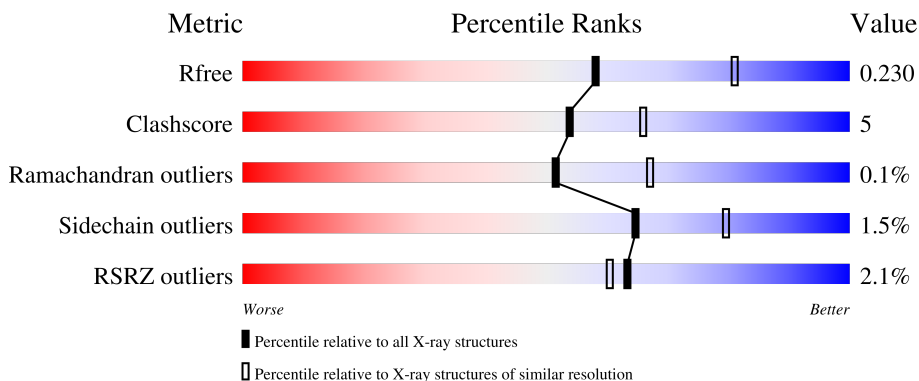|  |  |  |
|---|---|---|
| PDB ID | : | 6XUL |
| Title | : | Apo Ab 5b1 |
| Authors | : | Diskin, R.; Borenstein-Katz, A. |
| Deposited on | : | 2020-01-20 |
| Resolution | : | 2.41 Å(reported) |

# 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 2.41 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 | | 0.230 |
| Clashscore | | 5 |
| Ramachandran outliers | | 0.1% |
| Sidechain outliers | | 1.5% |
| 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 | 4647 (2.44-2.40) |
| Clashscore | 141614 | 5161 (2.44-2.40) |
| Ramachandran outliers | 138981 | 5073 (2.44-2.40) |
| Sidechain outliers | 138945 | 5074 (2.44-2.40) |
| RSRZ outliers | 127900 | 4543 (2.44-2.40) |

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 | 230 | .% / 83% 12% 5% |
| 1 | C | 230 | 3% / 84% 10% 5% |
| 1 | E | 230 | 8% / 85% 10% 5% |
| 1 | G | 230 | .% / 85% 10% 5% |
| 1 | H | 230 | .% / 83% 11% 5% |

*Continued on next page...*

*Continued from previous page...*

| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 1 | J | 230 | 2%<br>83%      11%    5% |
| 2 | B | 218 | %<br>89%      11% |
| 2 | D | 218 | %<br>87%      12% |
| 2 | F | 218 | 2%<br>88%      11% |
| 2 | I | 218 | 87%      13% |
| 2 | K | 218 | 2%<br>83%      16% |
| 2 | L | 218 | %<br>84%      14%   • |

# 2 Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 20926 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 Heavy chain.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|---------|-------|
| 1 | A | 218 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1665 | 1051 | 288 | 319 | 7 | | | |
| 1 | C | 218 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1665 | 1051 | 288 | 319 | 7 | | | |
| 1 | E | 218 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1665 | 1051 | 288 | 319 | 7 | | | |
| 1 | G | 219 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 1668 | 1052 | 289 | 321 | 6 | | | |
| 1 | J | 218 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1665 | 1051 | 288 | 319 | 7 | | | |
| 1 | H | 218 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1665 | 1051 | 288 | 319 | 7 | | | |

- Molecule 2 is a protein called Light chain.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|---------|-------|
| 2 | B | 217 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1651 | 1026 | 284 | 336 | 5 | | | |
| 2 | D | 217 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1651 | 1026 | 284 | 336 | 5 | | | |
| 2 | F | 217 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1651 | 1026 | 284 | 336 | 5 | | | |
| 2 | I | 217 | Total | C | N | O | S | 0 | 2 | 0 |
| | | | 1659 | 1031 | 287 | 336 | 5 | | | |
| 2 | K | 217 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 1648 | 1024 | 284 | 336 | 4 | | | |
| 2 | L | 217 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 1651 | 1026 | 284 | 336 | 5 | | | |

- Molecule 3 is water.

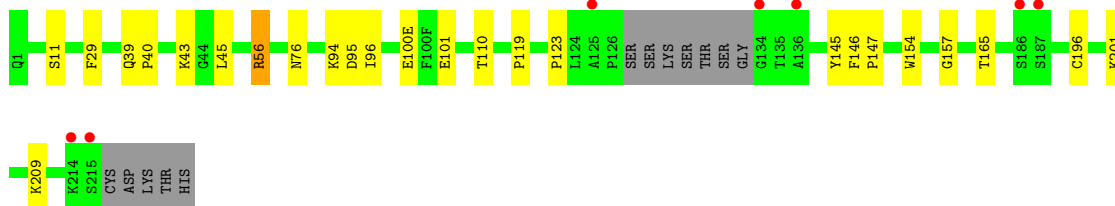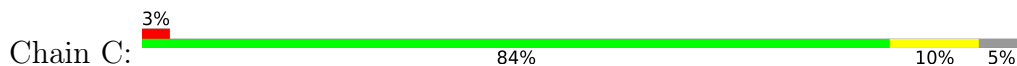| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|---|---|---|---|---|---|---|
| 3 | A | 138 | Total 138 | O 138 | 0 | 0 |
| 3 | B | 130 | Total 130 | O 130 | 0 | 0 |
| 3 | C | 103 | Total 103 | O 103 | 0 | 0 |
| 3 | D | 54 | Total 54 | O 54 | 0 | 0 |
| 3 | E | 98 | Total 98 | O 98 | 0 | 0 |
| 3 | F | 62 | Total 62 | O 62 | 0 | 0 |
| 3 | G | 125 | Total 125 | O 125 | 0 | 0 |
| 3 | I | 89 | Total 89 | O 89 | 0 | 0 |
| 3 | J | 85 | Total 85 | O 85 | 0 | 0 |
| 3 | K | 35 | Total 35 | O 35 | 0 | 0 |
| 3 | H | 62 | Total 62 | O 62 | 0 | 0 |
| 3 | L | 41 | Total 41 | O 41 | 0 | 0 |

# 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.
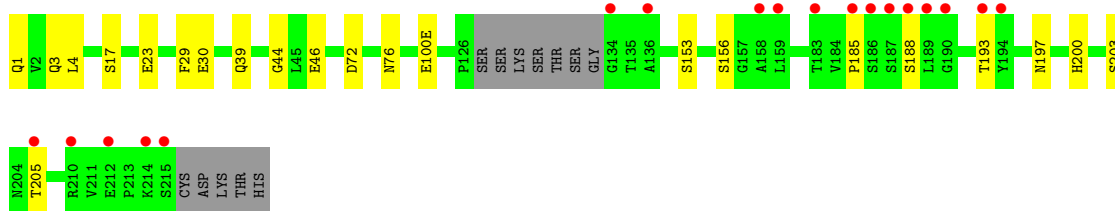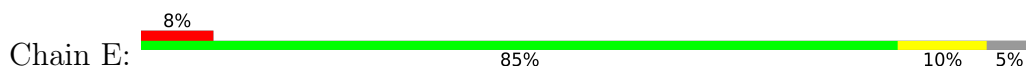
- Molecule 1: Heavy chain
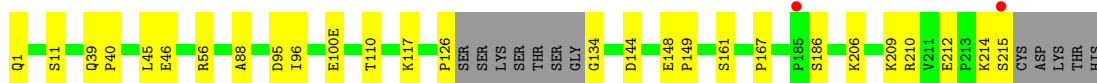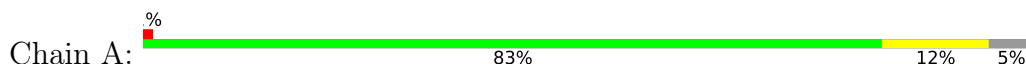
Chain A:



- Molecule 1: Heavy chain

Chain C:



- Molecule 1: Heavy chain

Chain E:



- Molecule 1: Heavy chain

Chain G:



- Molecule 1: Heavy chain

Chain J:



• Molecule 1: Heavy chain

Chain H:



• Molecule 2: Light chain

Chain B:



• Molecule 2: Light chain

Chain D:



• Molecule 2: Light chain

Chain F:



• Molecule 2: Light chain

Chain I:

● Molecule 2: Light chain

2%

Chain K:

83% 16%

Q1 L4 S9 R18 L39 R54 P55 S56 G57 V58 R61 S76 S80 A84 D92 D93 S94 L95 G95A G95B G99 V106 P113 P120 K126 V132 F139 R142 K145 V146 Q147 W148 E161 Y173 S176 S182 K183

A184 D185 K190 T197 S203 N210 R211 G212 E213 CYS

● Molecule 2: Light chain

%

Chain L:

84% 14% •

Q1 S2 T14 P15 S22 I30 W35 L39 L46 L47 I48 Y49 R54 P55 V58 P59 D60 R61 F62 S63 R66 S72 S76 E81 D82 E83 A84 D122 E123 D151 Q166 S171 K190 V191 L201 N210 E213 CYS
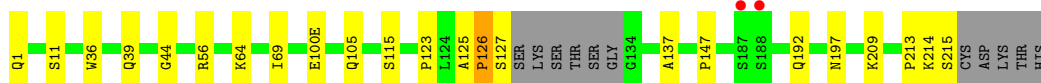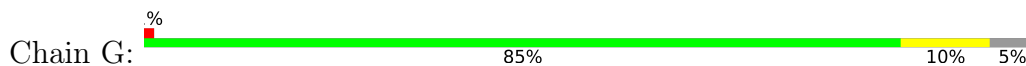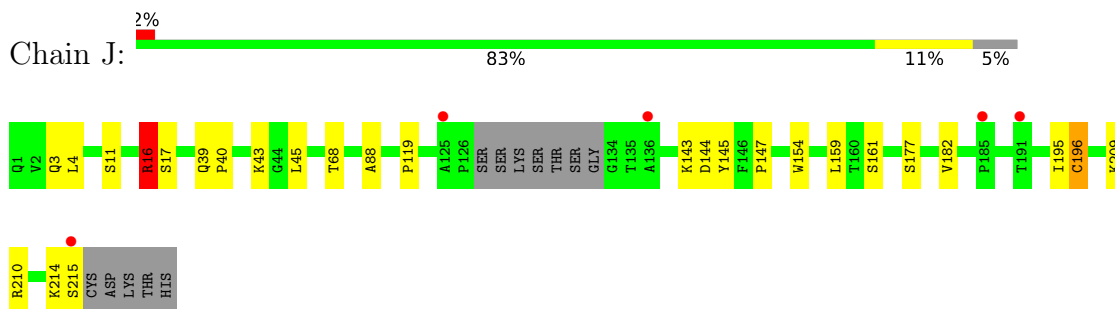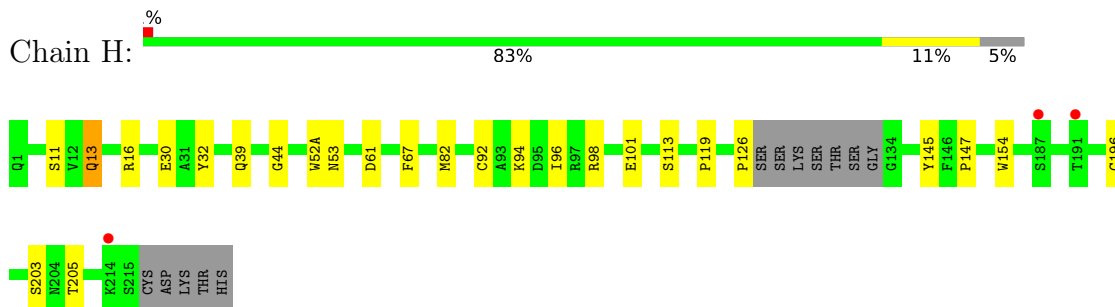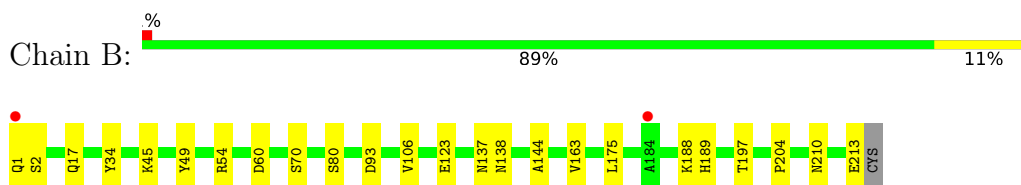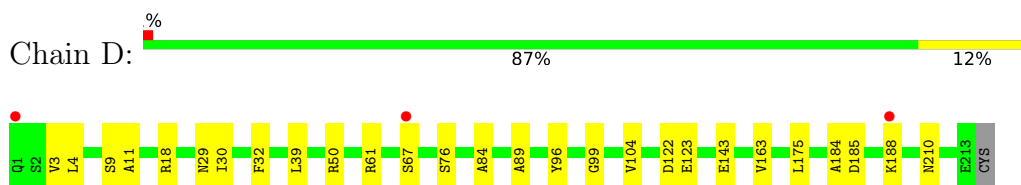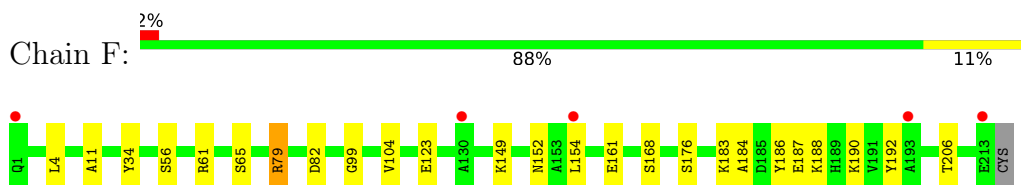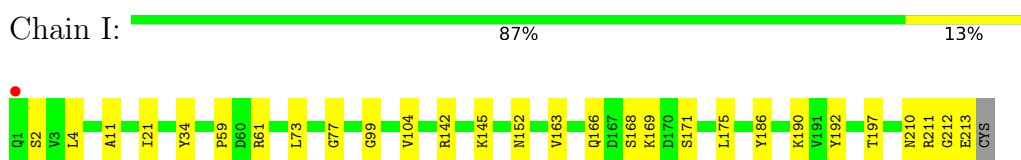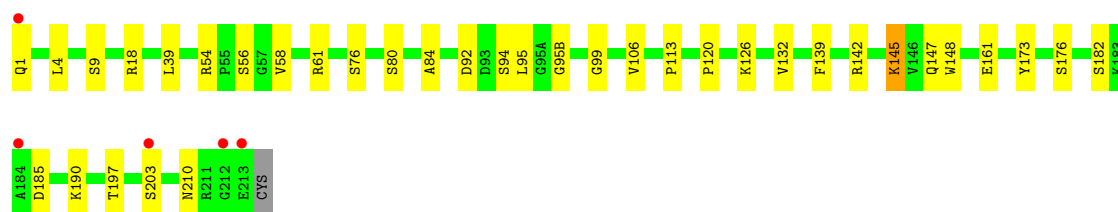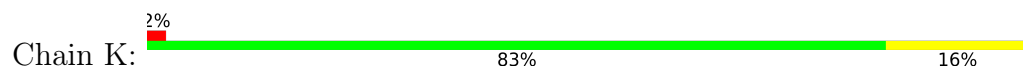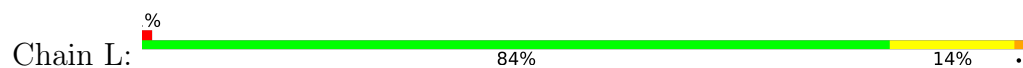
# 4 Data and refinement statistics ⓘ

| Property | Value | Source |
|:---:|:---:|:---:|
| Space group | P 32 | Depositor |
| Cell constants<br>a, b, c, $\alpha$, $\beta$, $\gamma$ | 155.02Å  155.02Å  121.78Å<br>90.00°  90.00°  120.00° | Depositor |
| Resolution (Å) | 47.88  –  2.41<br>47.88  –  2.41 | Depositor<br>EDS |
| % Data completeness<br>(in resolution range) | 99.7 (47.88-2.41)<br>99.7 (47.88-2.41) | Depositor<br>EDS |
| $R_{merge}$ | 0.15 | Depositor |
| $R_{sym}$ | (Not available) | Depositor |
| $< I/\sigma(I) >$ [1] | 1.36 (at 2.42Å) | Xtriage |
| Refinement program | PHENIX 1.16_3549 | Depositor |
| R, $R_{free}$ | 0.168  ,  0.230<br>0.168  ,  0.230 | Depositor<br>DCC |
| $R_{free}$ test set | 1976 reflections (1.57%) | wwPDB-VP |
| Wilson B-factor (Å$^2$) | 40.3 | Xtriage |
| Anisotropy | 0.358 | Xtriage |
| Bulk solvent $k_{sol}$(e/Å$^3$), $B_{sol}$(Å$^2$) | 0.31 , 44.3 | EDS |
| L-test for twinning[2] | $< |L| > = 0.48, < L^2 > = 0.31$ | Xtriage |
| Estimated twinning fraction | 0.026 for -h,-k,l<br>0.032 for h,-h-k,-l<br>0.017 for -k,-h,-l | Xtriage |
| $F_o$,$F_c$ correlation | 0.96 | EDS |
| Total number of atoms | 20926 | wwPDB-VP |
| Average B, all atoms (Å$^2$) | 49.0 | wwPDB-VP |

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 51.67 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.4292e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

---

[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 (i)

### 5.1 Standard geometry (i)

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 | 0.48 | 0/1709 | 0.63 | 0/2323 |
| 1 | C | 0.42 | 0/1709 | 0.60 | 0/2323 |
| 1 | E | 0.45 | 0/1709 | 0.61 | 0/2323 |
| 1 | G | 0.45 | 0/1709 | 0.61 | 0/2323 |
| 1 | H | 0.41 | 0/1709 | 0.59 | 0/2323 |
| 1 | J | 0.44 | 0/1709 | 0.59 | 1/2323 (0.0%) |
| 2 | B | 0.42 | 0/1691 | 0.60 | 0/2300 |
| 2 | D | 0.38 | 0/1691 | 0.57 | 0/2300 |
| 2 | F | 0.41 | 0/1691 | 0.59 | 0/2300 |
| 2 | I | 0.42 | 0/1702 | 0.59 | 0/2314 |
| 2 | K | 0.38 | 0/1685 | 0.53 | 0/2292 |
| 2 | L | 0.38 | 0/1691 | 0.57 | 0/2300 |
| All | All | 0.42 | 0/20405 | 0.59 | 1/27744 (0.0%) |

There are no bond length outliers.

All (1) bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed($^o$) | Ideal($^o$) |
|-----|-------|-----|------|-------|---|----------------|-------------|
| 1 | J | 16 | ARG | NE-CZ-NH2 | -5.03 | 117.79 | 120.30 |

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts (i)

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 | 1665 | 0 | 1624 | 21 | 0 |
| 1 | C | 1665 | 0 | 1624 | 19 | 0 |
| 1 | E | 1665 | 0 | 1624 | 16 | 0 |
| 1 | G | 1668 | 0 | 1625 | 17 | 0 |
| 1 | H | 1665 | 0 | 1624 | 19 | 0 |
| 1 | J | 1665 | 0 | 1624 | 17 | 0 |
| 2 | B | 1651 | 0 | 1588 | 17 | 0 |
| 2 | D | 1651 | 0 | 1588 | 17 | 0 |
| 2 | F | 1651 | 0 | 1588 | 16 | 0 |
| 2 | I | 1659 | 0 | 1601 | 15 | 0 |
| 2 | K | 1648 | 0 | 1584 | 20 | 0 |
| 2 | L | 1651 | 0 | 1588 | 19 | 0 |
| 3 | A | 138 | 0 | 0 | 7 | 0 |
| 3 | B | 130 | 0 | 0 | 6 | 0 |
| 3 | C | 103 | 0 | 0 | 5 | 0 |
| 3 | D | 54 | 0 | 0 | 4 | 0 |
| 3 | E | 98 | 0 | 0 | 7 | 1 |
| 3 | F | 62 | 0 | 0 | 4 | 0 |
| 3 | G | 125 | 0 | 0 | 9 | 1 |
| 3 | H | 62 | 0 | 0 | 3 | 0 |
| 3 | I | 89 | 0 | 0 | 2 | 0 |
| 3 | J | 85 | 0 | 0 | 4 | 0 |
| 3 | K | 35 | 0 | 0 | 6 | 0 |
| 3 | L | 41 | 0 | 0 | 6 | 0 |
| All | All | 20926 | 0 | 19282 | 207 | 1 |

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 5.

The worst 5 of 207 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) |
|--------|--------|--------------------------|-------------------|
| 1:E:193:THR:OG1 | 3:E:301:HOH:O | 1.85 | 0.92 |
| 2:F:168:SER:OG | 3:F:301:HOH:O | 1.88 | 0.91 |
| 2:L:201:LEU:O | 3:L:301:HOH:O | 1.88 | 0.90 |
| 1:C:43:LYS:O | 3:C:301:HOH:O | 1.93 | 0.86 |
| 2:B:144:ALA:O | 3:B:301:HOH:O | 1.95 | 0.85 |

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|---|---|---|---|
| 3:E:348:HOH:O | 3:G:421:HOH:O[2_545] | 2.08 | 0.12 |

## 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 | 215/230 (94%) | 211 (98%) | 4 (2%) | 0 | 100 | 100 |
| 1 | C | 215/230 (94%) | 207 (96%) | 8 (4%) | 0 | 100 | 100 |
| 1 | E | 215/230 (94%) | 205 (95%) | 10 (5%) | 0 | 100 | 100 |
| 1 | G | 215/230 (94%) | 209 (97%) | 5 (2%) | 1 (0%) | 29 | 40 |
| 1 | H | 215/230 (94%) | 209 (97%) | 6 (3%) | 0 | 100 | 100 |
| 1 | J | 215/230 (94%) | 209 (97%) | 6 (3%) | 0 | 100 | 100 |
| 2 | B | 216/218 (99%) | 211 (98%) | 5 (2%) | 0 | 100 | 100 |
| 2 | D | 216/218 (99%) | 207 (96%) | 9 (4%) | 0 | 100 | 100 |
| 2 | F | 216/218 (99%) | 208 (96%) | 8 (4%) | 0 | 100 | 100 |
| 2 | I | 217/218 (100%) | 209 (96%) | 6 (3%) | 2 (1%) | 17 | 24 |
| 2 | K | 215/218 (99%) | 207 (96%) | 8 (4%) | 0 | 100 | 100 |
| 2 | L | 216/218 (99%) | 205 (95%) | 11 (5%) | 0 | 100 | 100 |
| All | All | 2586/2688 (96%) | 2497 (97%) | 86 (3%) | 3 (0%) | 51 | 67 |

All (3) Ramachandran outliers are listed below:

| Mol | Chain | Res | Type |
|---|---|---|---|
| 1 | G | 126 | PRO |
| 2 | I | 77 | GLY |
| 2 | I | 212 | GLY |

### 5.3.2 Protein sidechains (i)

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 | 184/194 (95%) | 182 (99%) | 2 (1%) | 73 | 86 |
| 1 | C | 184/194 (95%) | 183 (100%) | 1 (0%) | 88 | 95 |
| 1 | E | 184/194 (95%) | 183 (100%) | 1 (0%) | 88 | 95 |
| 1 | G | 184/194 (95%) | 182 (99%) | 2 (1%) | 73 | 86 |
| 1 | H | 184/194 (95%) | 181 (98%) | 3 (2%) | 62 | 78 |
| 1 | J | 184/194 (95%) | 180 (98%) | 4 (2%) | 52 | 69 |
| 2 | B | 187/187 (100%) | 186 (100%) | 1 (0%) | 88 | 95 |
| 2 | D | 187/187 (100%) | 185 (99%) | 2 (1%) | 73 | 86 |
| 2 | F | 187/187 (100%) | 184 (98%) | 3 (2%) | 62 | 78 |
| 2 | I | 188/187 (100%) | 186 (99%) | 2 (1%) | 73 | 86 |
| 2 | K | 186/187 (100%) | 182 (98%) | 4 (2%) | 52 | 69 |
| 2 | L | 187/187 (100%) | 179 (96%) | 8 (4%) | 29 | 44 |
| All | All | 2226/2286 (97%) | 2193 (98%) | 33 (2%) | 65 | 79 |

5 of 33 residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 2 | L | 63 | SER |
| 2 | L | 72 | SER |
| 2 | L | 213 | GLU |
| 2 | I | 2 | SER |
| 1 | G | 197 | ASN |

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 2 | I | 137 | ASN |
| 2 | K | 210 | ASN |
| 2 | L | 95(C) | HIS |
| 2 | D | 189 | HIS |

*Continued on next page...*

*Continued from previous page...*

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 2 | B | 17 | GLN |

#### 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 ⓘ

There are no ligands in this entry.

## 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.

# 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 | 218/230 (94%) | -0.20 | 2 (0%) | 84 | 82 | 21, 34, 58, 94 | 0 |
| 1 | C | 218/230 (94%) | -0.06 | 7 (3%) | 47 | 45 | 26, 41, 84, 111 | 0 |
| 1 | E | 218/230 (94%) | 0.14 | 18 (8%) | 11 | 10 | 23, 43, 97, 121 | 0 |
| 1 | G | 219/230 (95%) | -0.19 | 2 (0%) | 84 | 82 | 24, 36, 64, 91 | 0 |
| 1 | H | 218/230 (94%) | -0.28 | 3 (1%) | 75 | 73 | 32, 49, 82, 106 | 0 |
| 1 | J | 218/230 (94%) | -0.22 | 5 (2%) | 60 | 57 | 31, 49, 84, 109 | 0 |
| 2 | B | 217/218 (99%) | -0.36 | 2 (0%) | 84 | 82 | 26, 38, 58, 91 | 0 |
| 2 | D | 217/218 (99%) | -0.15 | 3 (1%) | 75 | 73 | 33, 53, 85, 100 | 0 |
| 2 | F | 217/218 (99%) | -0.19 | 5 (2%) | 60 | 57 | 31, 50, 81, 118 | 0 |
| 2 | I | 217/218 (99%) | -0.40 | 1 (0%) | 91 | 89 | 29, 41, 58, 96 | 0 |
| 2 | K | 217/218 (99%) | -0.16 | 5 (2%) | 60 | 57 | 37, 63, 83, 102 | 0 |
| 2 | L | 217/218 (99%) | -0.27 | 2 (0%) | 84 | 82 | 41, 62, 81, 117 | 0 |
| All | All | 2611/2688 (97%) | -0.20 | 55 (2%) | 63 | 60 | 21, 46, 81, 121 | 0 |

The worst 5 of 55 RSRZ outliers are listed below:

| Mol | Chain | Res | Type | RSRZ |
|-----|-------|-----|------|------|
| 2 | L | 1 | GLN | 5.0 |
| 2 | B | 1 | GLN | 4.8 |
| 1 | J | 215 | SER | 4.7 |
| 2 | K | 212 | GLY | 4.5 |
| 1 | E | 189 | LEU | 4.3 |

## 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 ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.