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

Mar 20, 2024 – 12:05 PM EDT

| | | |
|---|---|---|
| PDB ID | : | 8EN8 |
| Title | : | Cross-reactive 3180 TCR recognition of HLA-B*35:01-NP4 epitope from 1972 influenza strain |
| Authors | : | Littler, D.R.; Rossjohn, J. |
| Deposited on | : | 2022-09-29 |
| Resolution | : | 2.70 Å(reported) |

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

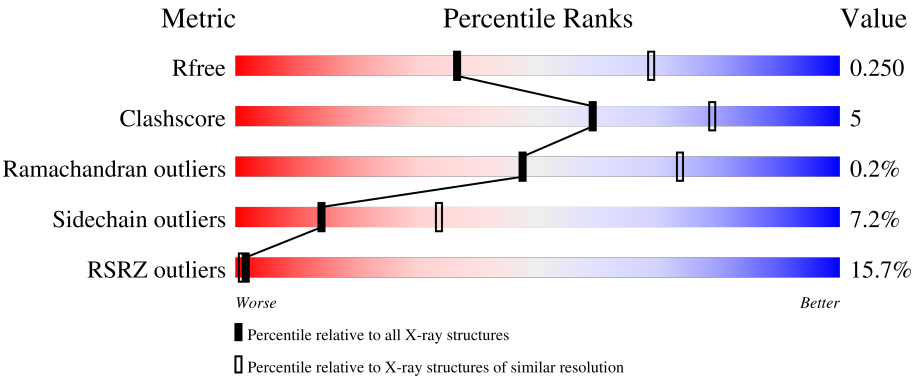
| | | |
|---|---|---|
| MolProbity | : | 4.02b-467 |
| Xtriage (Phenix) | : | 1.13 |
| EDS | : | 2.36 |
| Percentile statistics | : | 20191225.v01 (using entries in the PDB archive December 25th 2019) |
| Refmac | : | 5.8.0158 |
| CCP4 | : | 7.0.044 (Gargrove) |
| Ideal geometry (proteins) | : | Engh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

# 1 Overall quality at a glance (i)

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 | 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 | 276 | 17% 90% 8% |
| 1 | F | 276 | 14% 89% 10% |
| 2 | B | 100 | 11% 93% 5% |
| 2 | G | 100 | % 89% 9% |
| 3 | C | 9 | 78% 22% |

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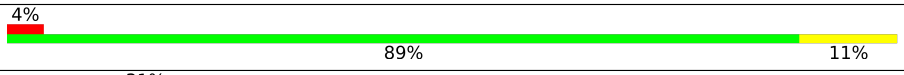
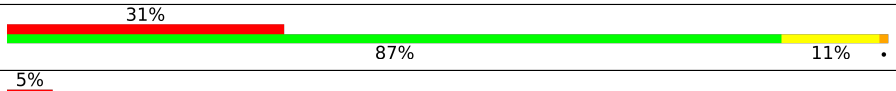
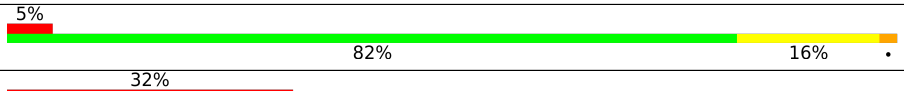
| Mol | Chain | Length | Quality of chain |
|---|---|---|---|
| 3 | H | 9 |   |
| 4 | D | 205 |   |
| 4 | I | 205 |   |
| 5 | E | 245 |   |
| 5 | J | 245 |   |

# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 13859 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 MHC class I antigen.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|---|---|---|---|---|---|---|---|---|---|---|
| 1 | A | 276 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 2262 | 1409 | 413 | 433 | 7 | | | |
| 1 | F | 276 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 2262 | 1409 | 413 | 433 | 7 | | | |

- Molecule 2 is a protein called Beta-2-microglobulin.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|---|---|---|---|---|---|---|---|---|---|---|
| 2 | B | 99 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 829 | 528 | 140 | 158 | 3 | | | |
| 2 | G | 99 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 829 | 528 | 140 | 158 | 3 | | | |

There are 2 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|---|---|---|---|---|---|
| B | 0 | MET | - | initiating methionine | UNP P61769 |
| G | 0 | MET | - | initiating methionine | UNP P61769 |

- Molecule 3 is a protein called Nucleoprotein NP4 epitope.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|---|---|---|---|---|---|---|---|---|---|---|
| 3 | C | 9 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 73 | 48 | 10 | 14 | 1 | | | |
| 3 | H | 9 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 73 | 48 | 10 | 14 | 1 | | | |

- Molecule 4 is a protein called 3180 TCR alpha chain.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|---|---|---|---|---|---|---|---|---|---|---|
| 4 | D | 205 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 1594 | 994 | 266 | 326 | 8 | | | |

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| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|---|---|---|---|---|---|---|---|---|---|---|
| 4 | I | 205 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 1594 | 994 | 266 | 326 | 8 | | | |

- Molecule 5 is a protein called 3180 TCR beta chain.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|---|---|---|---|---|---|---|---|---|---|---|
| 5 | E | 245 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 1937 | 1219 | 334 | 375 | 9 | | | |
| 5 | J | 245 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 1937 | 1219 | 334 | 375 | 9 | | | |

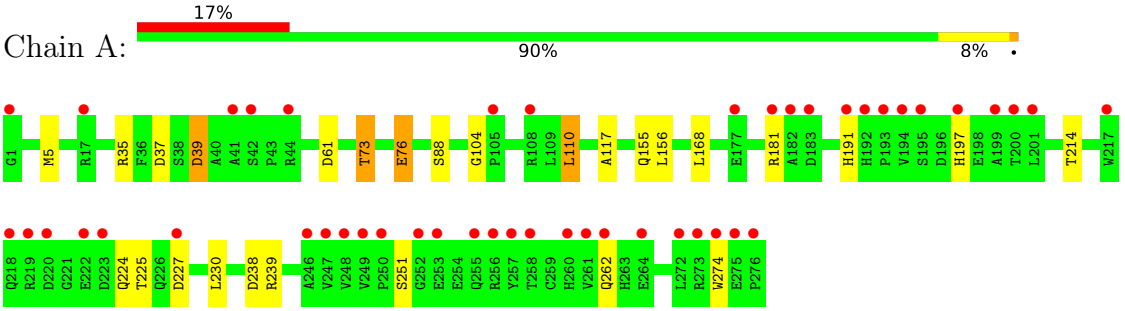
- Molecule 6 is water.

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|---|---|---|---|---|---|---|
| 6 | A | 79 | Total | O | 0 | 0 |
| | | | 79 | 79 | | |
| 6 | B | 14 | Total | O | 0 | 0 |
| | | | 14 | 14 | | |
| 6 | C | 3 | Total | O | 0 | 0 |
| | | | 3 | 3 | | |
| 6 | D | 51 | Total | O | 0 | 0 |
| | | | 51 | 51 | | |
| 6 | E | 80 | Total | O | 0 | 0 |
| | | | 80 | 80 | | |
| 6 | F | 110 | Total | O | 0 | 0 |
| | | | 110 | 110 | | |
| 6 | G | 36 | Total | O | 0 | 0 |
| | | | 36 | 36 | | |
| 6 | H | 8 | Total | O | 0 | 0 |
| | | | 8 | 8 | | |
| 6 | I | 34 | Total | O | 0 | 0 |
| | | | 34 | 34 | | |
| 6 | J | 54 | Total | O | 0 | 0 |
| | | | 54 | 54 | | |

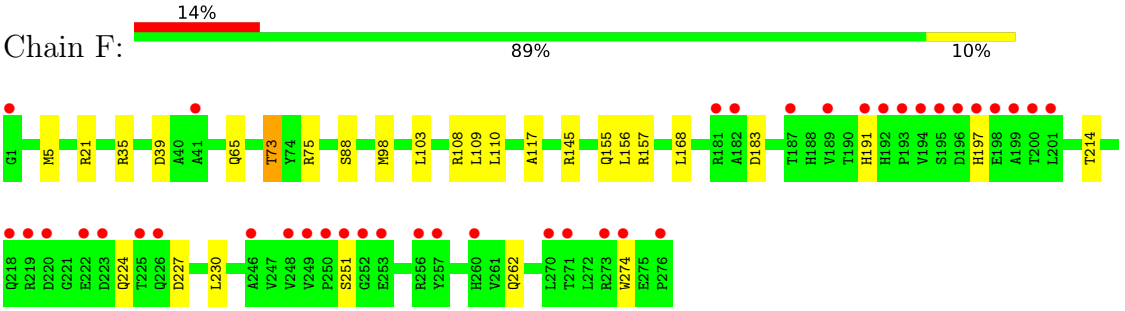
# 3 Residue-property plots (i)

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.

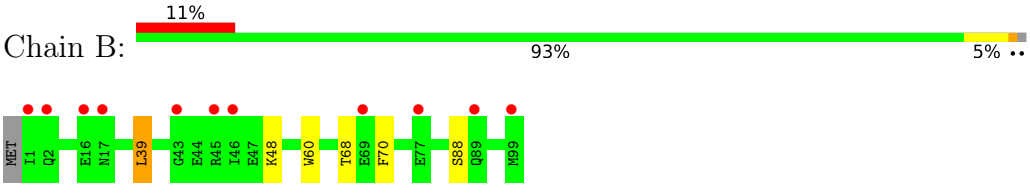
• Molecule 1: MHC class I antigen



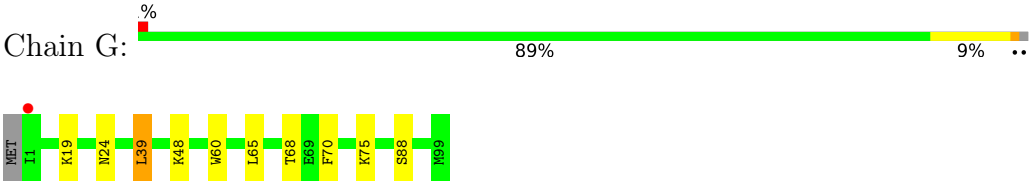

• Molecule 1: MHC class I antigen




• Molecule 2: Beta-2-microglobulin

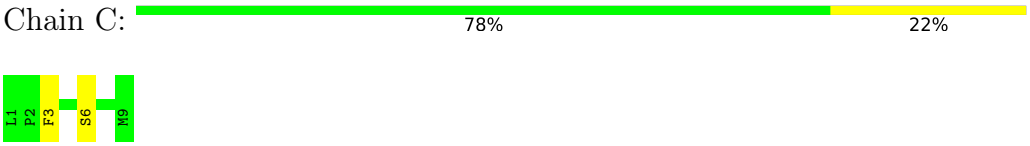



• Molecule 2: Beta-2-microglobulin



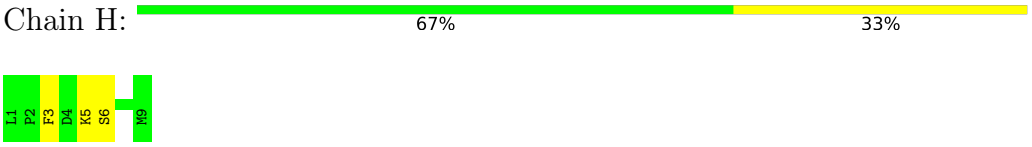

• Molecule 3: Nucleoprotein NP4 epitope

Chain C:



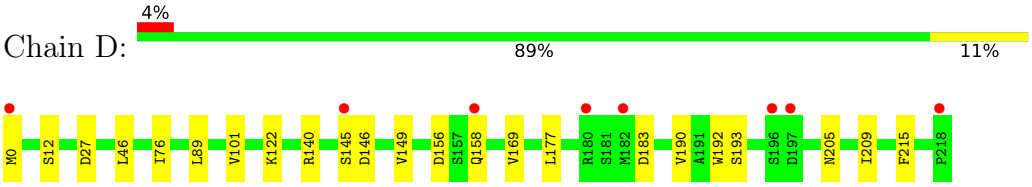

• Molecule 3: Nucleoprotein NP4 epitope

Chain H:



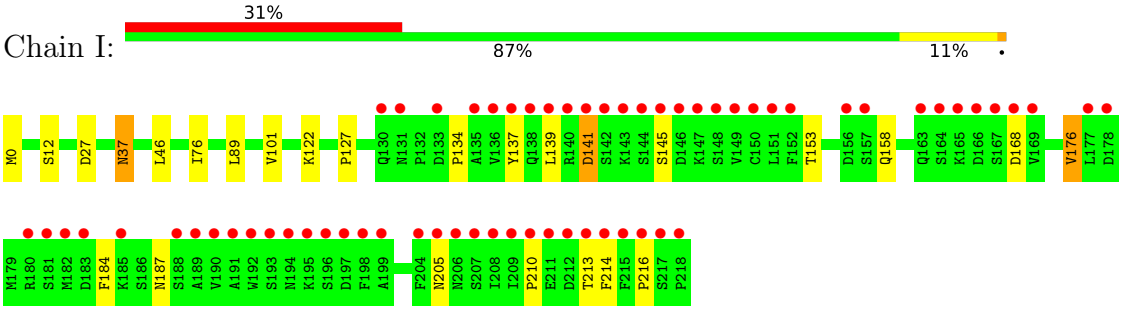

• Molecule 4: 3180 TCR alpha chain

Chain D:



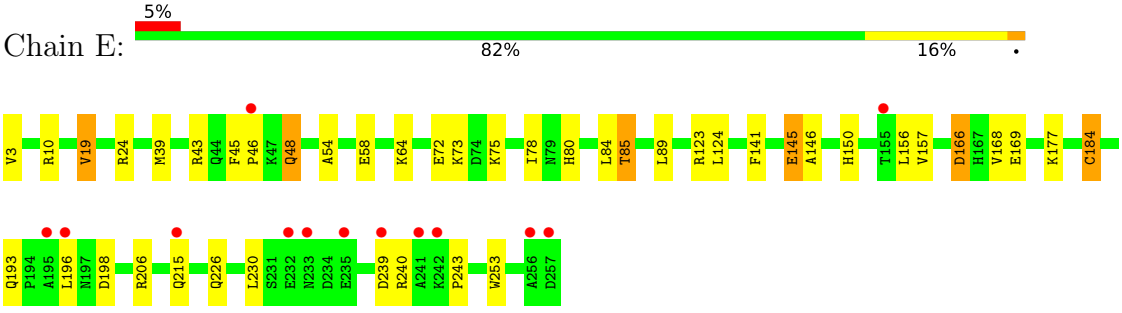

• Molecule 4: 3180 TCR alpha chain

Chain I:



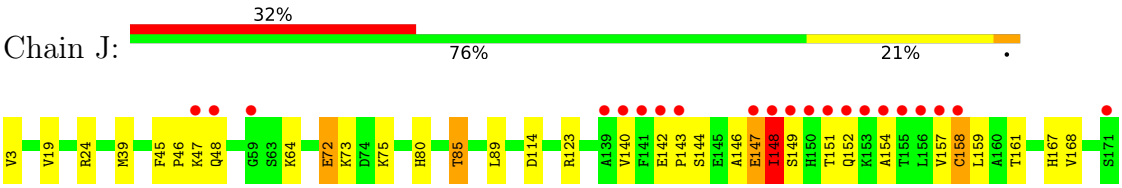

• Molecule 5: 3180 TCR beta chain

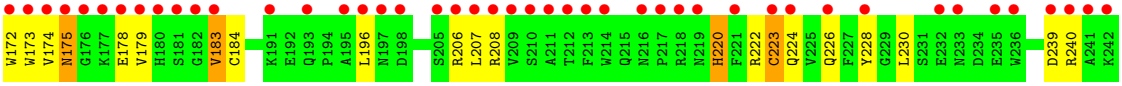
Chain E:

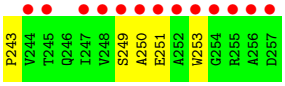



• Molecule 5: 3180 TCR beta chain

Chain J:




W172
W173
V174
N175
G176
K177
E178
V179
H180
S181
G182
V183
C184
K191
E192
Q193
P194
A195
L196
N197
D198
S205
R206
L207
R208
V209
S210
A211
T212
F213
W214
Q215
N216
P217
R218
N219
H220
F221
R222
C223
Q224
V225
Q226
F227
Y228
G229
L230
S231
E232
N233
D234
E235
W236
D239
R240
A241
K242


P243
V244
T245
Q246
I247
V248
S249
A250
E251
A252
W253
G254
R255
A256
D257

# 4 Data and refinement statistics (i)

| Property | Value | Source |
|---|---|---|
| Space group | P 21 21 21 | Depositor |
| Cell constants<br>a, b, c, $\alpha$, $\beta$, $\gamma$ | 79.77Å 116.69Å 253.18Å<br>90.00° 90.00° 90.00° | Depositor |
| Resolution (Å) | 27.20 – 2.70<br>27.20 – 2.70 | Depositor<br>EDS |
| % Data completeness<br>(in resolution range) | 99.9 (27.20-2.70)<br>100.0 (27.20-2.70) | Depositor<br>EDS |
| $R_{merge}$ | (Not available) | Depositor |
| $R_{sym}$ | (Not available) | Depositor |
| $< I/\sigma(I) >$ [1] | 2.12 (at 2.72Å) | Xtriage |
| Refinement program | BUSTER, PHENIX 1.10.1_2155 | Depositor |
| R, $R_{free}$ | 0.207 , 0.249<br>0.211 , 0.250 | Depositor<br>DCC |
| $R_{free}$ test set | 3281 reflections (4.99%) | wwPDB-VP |
| Wilson B-factor (Å$^2$) | 46.2 | Xtriage |
| Anisotropy | 0.458 | Xtriage |
| Bulk solvent $k_{sol}$(e/Å$^3$), $B_{sol}$(Å$^2$) | 0.30 , 61.9 | EDS |
| L-test for twinning[2] | $< \|L\| > = 0.48$, $< L^2 > = 0.31$ | Xtriage |
| Estimated twinning fraction | No twinning to report. | Xtriage |
| $F_o$,$F_c$ correlation | 0.92 | EDS |
| Total number of atoms | 13859 | wwPDB-VP |
| Average B, all atoms (Å$^2$) | 83.0 | wwPDB-VP |

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% of the height of the origin peak. No significant pseudotranslation is detected.*

[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

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 | 0.45 | 0/2325 | 0.65 | 0/3161 |
| 1 | F | 0.47 | 0/2325 | 0.64 | 0/3161 |
| 2 | B | 0.43 | 0/852 | 0.62 | 0/1152 |
| 2 | G | 0.42 | 0/852 | 0.64 | 0/1152 |
| 3 | C | 0.67 | 0/74 | 0.68 | 0/97 |
| 3 | H | 0.40 | 0/74 | 0.67 | 0/97 |
| 4 | D | 0.41 | 0/1628 | 0.64 | 0/2205 |
| 4 | I | 0.42 | 0/1628 | 0.63 | 0/2205 |
| 5 | E | 0.40 | 0/1988 | 0.68 | 0/2705 |
| 5 | J | 0.45 | 0/1988 | 0.70 | 0/2705 |
| All | All | 0.44 | 0/13734 | 0.66 | 0/18640 |

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 | 2262 | 0 | 2122 | 13 | 0 |
| 1 | F | 2262 | 0 | 2122 | 10 | 0 |
| 2 | B | 829 | 0 | 794 | 2 | 0 |
| 2 | G | 829 | 0 | 794 | 3 | 0 |
| 3 | C | 73 | 0 | 78 | 2 | 0 |
| 3 | H | 73 | 0 | 78 | 3 | 0 |

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| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|---|---|---|---|---|---|---|
| 4 | D | 1594 | 0 | 1521 | 7 | 0 |
| 4 | I | 1594 | 0 | 1521 | 23 | 0 |
| 5 | E | 1937 | 0 | 1856 | 30 | 0 |
| 5 | J | 1937 | 0 | 1856 | 57 | 0 |
| 6 | A | 79 | 0 | 0 | 0 | 0 |
| 6 | B | 14 | 0 | 0 | 0 | 0 |
| 6 | C | 3 | 0 | 0 | 0 | 0 |
| 6 | D | 51 | 0 | 0 | 0 | 0 |
| 6 | E | 80 | 0 | 0 | 2 | 0 |
| 6 | F | 110 | 0 | 0 | 1 | 0 |
| 6 | G | 36 | 0 | 0 | 0 | 0 |
| 6 | H | 8 | 0 | 0 | 0 | 0 |
| 6 | I | 34 | 0 | 0 | 1 | 0 |
| 6 | J | 54 | 0 | 0 | 0 | 0 |
| All | All | 13859 | 0 | 12742 | 124 | 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 5.

The worst 5 of 124 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) |
|---|---|---|---|
| 5:J:147:GLU:O | 5:J:148:ILE:CG1 | 1.95 | 1.14 |
| 5:J:147:GLU:O | 5:J:148:ILE:HG12 | 1.52 | 1.08 |
| 4:I:127:PRO:HG3 | 4:I:176:VAL:HG11 | 1.40 | 1.03 |
| 5:J:175:ASN:H | 5:J:220:HIS:HD2 | 1.08 | 1.00 |
| 5:J:147:GLU:C | 5:J:148:ILE:HG12 | 1.77 | 1.00 |

There are no symmetry-related clashes.

## 5.3 Torsion angles (i)

### 5.3.1 Protein backbone (i)

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 | 275/276 (100%) | 265 (96%) | 9 (3%) | 1 (0%) | 34 | 60 |
| 1 | F | 275/276 (100%) | 262 (95%) | 12 (4%) | 1 (0%) | 34 | 60 |
| 2 | B | 97/100 (97%) | 95 (98%) | 2 (2%) | 0 | 100 | 100 |
| 2 | G | 97/100 (97%) | 96 (99%) | 1 (1%) | 0 | 100 | 100 |
| 3 | C | 7/9 (78%) | 7 (100%) | 0 | 0 | 100 | 100 |
| 3 | H | 7/9 (78%) | 7 (100%) | 0 | 0 | 100 | 100 |
| 4 | D | 203/205 (99%) | 195 (96%) | 8 (4%) | 0 | 100 | 100 |
| 4 | I | 203/205 (99%) | 191 (94%) | 12 (6%) | 0 | 100 | 100 |
| 5 | E | 243/245 (99%) | 232 (96%) | 10 (4%) | 1 (0%) | 34 | 60 |
| 5 | J | 243/245 (99%) | 224 (92%) | 18 (7%) | 1 (0%) | 34 | 60 |
| All | All | 1650/1670 (99%) | 1574 (95%) | 72 (4%) | 4 (0%) | 47 | 73 |

All (4) Ramachandran outliers are listed below:

| Mol | Chain | Res | Type |
|---|---|---|---|
| 5 | J | 148 | ILE |
| 1 | A | 251 | SER |
| 5 | E | 166 | ASP |
| 1 | F | 251 | SER |

### 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 | 235/234 (100%) | 224 (95%) | 11 (5%) | 26 | 54 |
| 1 | F | 235/234 (100%) | 222 (94%) | 13 (6%) | 21 | 46 |
| 2 | B | 94/95 (99%) | 90 (96%) | 4 (4%) | 29 | 57 |
| 2 | G | 94/95 (99%) | 88 (94%) | 6 (6%) | 17 | 39 |
| 3 | C | 9/9 (100%) | 9 (100%) | 0 | 100 | 100 |
| 3 | H | 9/9 (100%) | 9 (100%) | 0 | 100 | 100 |
| 4 | D | 182/182 (100%) | 169 (93%) | 13 (7%) | 14 | 34 |

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| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles |
|---|---|---|---|---|---|
| 4 | I | 182/182 (100%) | 170 (93%) | 12 (7%) | 16 38 |
| 5 | E | 216/216 (100%) | 196 (91%) | 20 (9%) | 9 21 |
| 5 | J | 216/216 (100%) | 189 (88%) | 27 (12%) | 4 10 |
| All | All | 1472/1472 (100%) | 1366 (93%) | 106 (7%) | 14 34 |

5 of 106 residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|---|---|---|---|
| 1 | F | 157 | ARG |
| 4 | I | 46 | LEU |
| 5 | J | 196 | LEU |
| 1 | F | 197 | HIS |
| 2 | G | 75 | LYS |

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

| Mol | Chain | Res | Type |
|---|---|---|---|
| 5 | J | 175 | ASN |
| 5 | J | 220 | HIS |
| 5 | J | 224 | GLN |
| 5 | E | 48 | GLN |
| 4 | D | 202 | ASN |

### 5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry (i)

There are no ligands in this entry.

## 5.7 Other polymers (i)

There are no such residues in this entry.

## 5.8 Polymer linkage issues (i)

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 | 276/276 (100%) | 0.63 | 47 (17%) 1 1 | 29, 68, 177, 207 | 7 (2%) |
| 1 | F | 276/276 (100%) | 0.36 | 39 (14%) 2 1 | 21, 52, 185, 205 | 8 (2%) |
| 2 | B | 99/100 (99%) | 0.65 | 11 (11%) 5 4 | 47, 90, 127, 138 | 6 (6%) |
| 2 | G | 99/100 (99%) | 0.00 | 1 (1%) 82 83 | 32, 62, 100, 115 | 6 (6%) |
| 3 | C | 9/9 (100%) | 0.09 | 0 100 100 | 29, 34, 43, 50 | 0 |
| 3 | H | 9/9 (100%) | 0.25 | 0 100 100 | 26, 32, 36, 39 | 0 |
| 4 | D | 205/205 (100%) | 0.06 | 8 (3%) 39 38 | 32, 61, 105, 139 | 1 (0%) |
| 4 | I | 205/205 (100%) | 1.88 | 64 (31%) 0 0 | 32, 91, 226, 253 | 0 |
| 5 | E | 245/245 (100%) | 0.12 | 13 (5%) 26 25 | 27, 57, 103, 144 | 0 |
| 5 | J | 245/245 (100%) | 1.89 | 79 (32%) 0 0 | 27, 107, 230, 263 | 0 |
| All | All | 1668/1670 (99%) | 0.74 | 262 (15%) 2 1 | 21, 67, 193, 263 | 28 (1%) |

The worst 5 of 262 RSRZ outliers are listed below:

| Mol | Chain | Res | Type | RSRZ |
|---|---|---|---|---|
| 5 | J | 257 | ASP | 21.5 |
| 4 | I | 206 | ASN | 17.5 |
| 5 | J | 256 | ALA | 16.4 |
| 4 | I | 218 | PRO | 14.1 |
| 5 | J | 216 | ASN | 14.0 |

## 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 (i)

There are no monosaccharides in this entry.

## 6.4 Ligands (i)

There are no ligands in this entry.

## 6.5 Other polymers (i)

There are no such residues in this entry.