



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 5BYC
Title : Crystal structure of human ribokinase in C2 spacegroup
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Deposited on : 2015-06-10
Resolution : 1.95 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

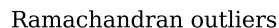
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027674
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027674

i

X-RAY DIFFRACTION

A.

 R_{free} 1

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5254 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 Ribokinase.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|---------|-------|
| 1 | A | 318 | Total | C | N | O | S | 0 | 3 | 0 |
| | | | 2393 | 1510 | 398 | 471 | 14 | | | |
| 1 | B | 316 | Total | C | N | O | S | 0 | 2 | 0 |
| | | | 2347 | 1482 | 391 | 460 | 14 | | | |

There are 16 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------------|------------|
| A | 323 | LEU | - | expression tag | UNP Q9H477 |
| A | 324 | GLU | - | expression tag | UNP Q9H477 |
| A | 325 | HIS | - | expression tag | UNP Q9H477 |
| A | 326 | HIS | - | expression tag | UNP Q9H477 |
| A | 327 | HIS | - | expression tag | UNP Q9H477 |
| A | 328 | HIS | - | expression tag | UNP Q9H477 |
| A | 329 | HIS | - | expression tag | UNP Q9H477 |
| A | 330 | HIS | - | expression tag | UNP Q9H477 |
| B | 323 | LEU | - | expression tag | UNP Q9H477 |
| B | 324 | GLU | - | expression tag | UNP Q9H477 |
| B | 325 | HIS | - | expression tag | UNP Q9H477 |
| B | 326 | HIS | - | expression tag | UNP Q9H477 |
| B | 327 | HIS | - | expression tag | UNP Q9H477 |
| B | 328 | HIS | - | expression tag | UNP Q9H477 |
| B | 329 | HIS | - | expression tag | UNP Q9H477 |
| B | 330 | HIS | - | expression tag | UNP Q9H477 |

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---------|---------|
| 2 | B | 1 | Total | Na | 0 | 0 |
| | | | 1 | 1 | | |
| 2 | A | 1 | Total | Na | 0 | 0 |
| | | | 1 | 1 | | |

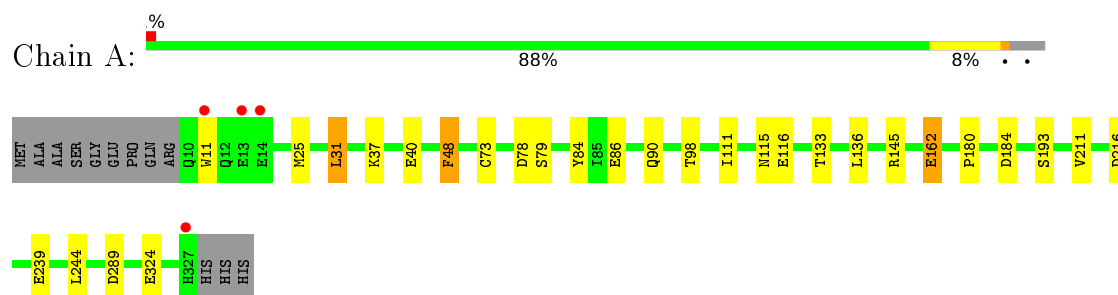
- Molecule 3 is water.

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|--------------|----------|---------|---------|
| 3 | A | 254 | Total 255 | O 255 | 0 | 1 |
| 3 | B | 257 | Total 257 | O 257 | 0 | 0 |

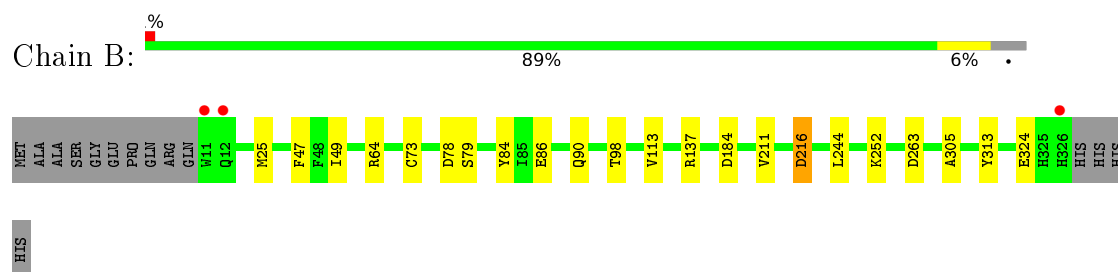
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Ribokinase



• Molecule 1: Ribokinase



4 Data and refinement statistics

| Property | Value | Source |
|---|---|------------------|
| Space group | C 1 2 1 | Depositor |
| Cell constants a, b, c, α , β , γ | 180.92Å 47.04Å 80.91Å 90.00° 98.73° 90.00° | Depositor |
| Resolution (Å) | 45.49 – 1.95 45.49 – 1.95 | Depositor EDS |
| % Data completeness (in resolution range) | 99.8 (45.49-1.95) 99.8 (45.49-1.95) | Depositor EDS |
| R_{merge} | 0.13 | Depositor |
| R_{sym} | (Not available) | Depositor |
| $\langle I/\sigma(I) \rangle$ ¹ | 1.99 (at 1.95Å) | Xtriage |
| Refinement program | REFMAC 5.8.0123 | Depositor |
| R, R_{free} | 0.173 , 0.220 0.183 , 0.227 | Depositor DCC |
| R_{free} test set | 2508 reflections (5.35%) | DCC |
| Wilson B-factor (Å ²) | 22.6 | Xtriage |
| Anisotropy | 0.300 | Xtriage |
| Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²) | 0.33 , 39.1 | EDS |
| L-test for twinning ² | $\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$ | Xtriage |
| Estimated twinning fraction | No twinning to report. | Xtriage |
| F_o, F_c correlation | 0.96 | EDS |
| Total number of atoms | 5254 | wwPDB-VP |
| Average B, all atoms (Å ²) | 28.0 | wwPDB-VP |

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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 | 1.20 | 8/2431 (0.3%) | 1.08 | 9/3306 (0.3%) |
| 1 | B | 1.14 | 3/2384 (0.1%) | 1.08 | 7/3246 (0.2%) |
| All | All | 1.17 | 11/4815 (0.2%) | 1.08 | 16/6552 (0.2%) |

The worst 5 of 11 bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|--------|-------|-------------|----------|
| 1 | B | 79 | SER | CB-OG | -9.30 | 1.30 | 1.42 |
| 1 | B | 324 | GLU | CD-OE2 | 6.91 | 1.33 | 1.25 |
| 1 | A | 40 | GLU | CD-OE1 | -6.75 | 1.18 | 1.25 |
| 1 | A | 324 | GLU | CD-OE2 | 6.49 | 1.32 | 1.25 |
| 1 | A | 48 | PHE | CG-CD1 | 5.83 | 1.47 | 1.38 |

The worst 5 of 16 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|-----|------|-----------|--------|-------------|----------|
| 1 | B | 25 | MET | CG-SD-CE | -10.54 | 83.34 | 100.20 |
| 1 | A | 25 | MET | CG-SD-CE | -10.27 | 83.77 | 100.20 |
| 1 | B | 137 | ARG | NE-CZ-NH1 | 7.49 | 124.04 | 120.30 |
| 1 | B | 244 | LEU | CA-CB-CG | 7.06 | 131.54 | 115.30 |
| 1 | A | 145 | ARG | NE-CZ-NH2 | -6.71 | 116.95 | 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 | 2393 | 0 | 2383 | 16 | 0 |
| 1 | B | 2347 | 0 | 2333 | 6 | 0 |
| 2 | A | 1 | 0 | 0 | 0 | 0 |
| 2 | B | 1 | 0 | 0 | 0 | 0 |
| 3 | A | 255 | 0 | 0 | 4 | 0 |
| 3 | B | 257 | 0 | 0 | 0 | 0 |
| All | All | 5254 | 0 | 4716 | 22 | 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 2.

The worst 5 of 22 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:A:31[A]:LEU:HG | 1:A:48:PHE:HE2 | 1.45 | 0.79 |
| 1:A:133:THR:HG23 | 1:A:162:GLU:OE2 | 1.90 | 0.70 |
| 1:A:31[A]:LEU:CG | 1:A:48:PHE:HE2 | 2.11 | 0.64 |
| 1:A:111:ILE:HD11 | 3:A:505:HOH:O | 2.02 | 0.58 |
| 1:A:115:ASN:HB2 | 3:A:519:HOH:O | 2.09 | 0.53 |

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 | 319/330 (97%) | 309 (97%) | 10 (3%) | 0 | 100 | 100 |
| 1 | B | 316/330 (96%) | 309 (98%) | 7 (2%) | 0 | 100 | 100 |
| All | All | 635/660 (96%) | 618 (97%) | 17 (3%) | 0 | 100 | 100 |

There are no Ramachandran outliers to report.

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 | 258/266 (97%) | 255 (99%) | 3 (1%) | 78 | 75 |
| 1 | B | 251/266 (94%) | 249 (99%) | 2 (1%) | 86 | 85 |
| All | All | 509/532 (96%) | 504 (99%) | 5 (1%) | 82 | 80 |

All (5) residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | A | 84 | TYR |
| 1 | A | 180 | PRO |
| 1 | A | 244 | LEU |
| 1 | B | 84 | TYR |
| 1 | B | 113 | VAL |

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | A | 43 | HIS |
| 1 | A | 325 | HIS |
| 1 | B | 90 | 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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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, 95th 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 | 318/330 (96%) | -0.27 | 4 (1%) 79 86 | 13, 25, 48, 77 | 0 |
| 1 | B | 316/330 (95%) | -0.30 | 3 (0%) 85 90 | 14, 27, 52, 82 | 0 |
| All | All | 634/660 (96%) | -0.29 | 7 (1%) 82 88 | 13, 26, 49, 82 | 0 |

The worst 5 of 7 RSRZ outliers are listed below:

| Mol | Chain | Res | Type | RSRZ |
|-----|-------|-----|------|------|
| 1 | A | 327 | HIS | 3.3 |
| 1 | A | 11 | TRP | 3.0 |
| 1 | B | 326 | HIS | 2.7 |
| 1 | B | 12 | GLN | 2.6 |
| 1 | A | 14 | GLU | 2.5 |

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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 | LLDF | B-factors(\AA^2) | Q<0.9 |
|-----|------|-------|-----|-------|------|------|-------|-----------------------------|-------|
| 2 | NA | B | 401 | 1/1 | 0.97 | 0.07 | -0.64 | 24,24,24,24 | 0 |
| 2 | NA | A | 401 | 1/1 | 0.98 | 0.06 | -1.14 | 24,24,24,24 | 0 |

6.5 Other polymers [i](#)

There are no such residues in this entry.