



Full wwPDB X-ray Structure Validation Report i

May 21, 2020 – 02:44 am BST

PDB ID : 4O4U
Title : Crystal structure of the vaccine antigen Transferrin Binding Protein B (TbpB) mutant Trp-176-Ala from Haemophilus parasuis Hp5
Authors : Calmettes, C.; Yu, R.H.; Schryvers, A.B.; Moraes, T.F.
Deposited on : 2013-12-19
Resolution : 2.64 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 following versions of software and data (see [references](#) ①) were used in the production of this report:

| | | |
|--------------------------------|---|--|
| MolProbity | : | 4.02b-467 |
| Mogul | : | 1.8.5 (274361), CSD as541be (2020) |
| Xtriage (Phenix) | : | 1.13 |
| EDS | : | 2.11 |
| 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) | : | 2.11 |

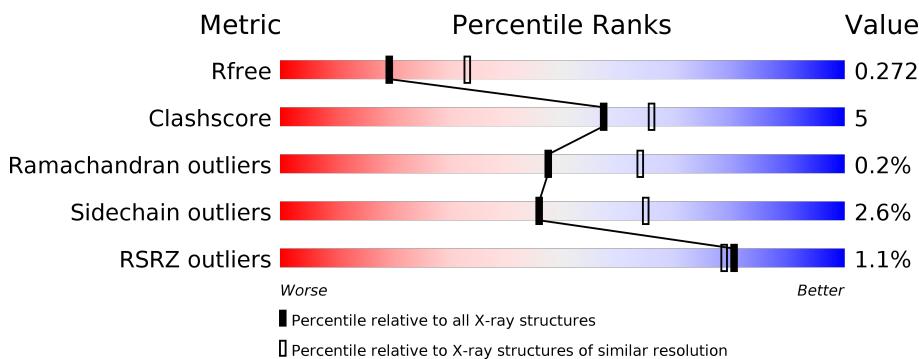
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

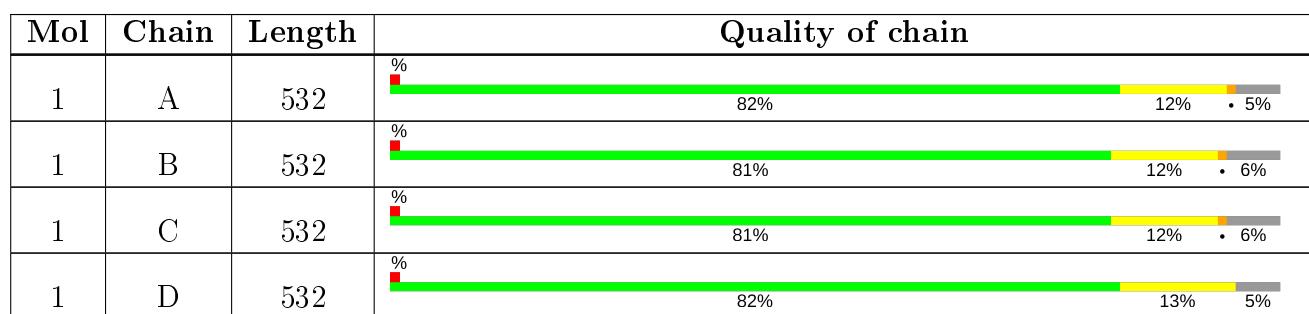
The reported resolution of this entry is 2.64 Å.

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 | 1426 (2.66-2.62) |
| Clashscore | 141614 | 1472 (2.66-2.62) |
| Ramachandran outliers | 138981 | 1446 (2.66-2.62) |
| Sidechain outliers | 138945 | 1446 (2.66-2.62) |
| RSRZ outliers | 127900 | 1408 (2.66-2.62) |

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 on 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.



2 Entry composition [\(i\)](#)

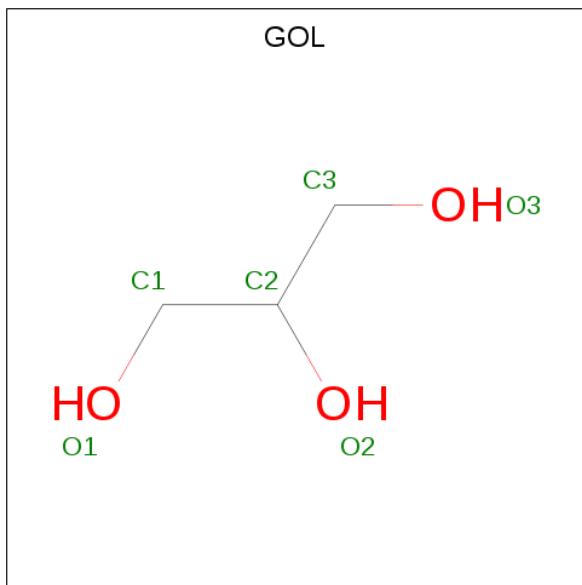
There are 3 unique types of molecules in this entry. The entry contains 15892 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 TbpB.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|---------|-------|
| 1 | A | 505 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 3912 | 2465 | 655 | 782 | 10 | | | |
| 1 | B | 500 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 3871 | 2441 | 644 | 776 | 10 | | | |
| 1 | C | 499 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 3861 | 2434 | 643 | 774 | 10 | | | |
| 1 | D | 507 | Total | C | N | O | S | 0 | 2 | 0 |
| | | | 3944 | 2487 | 659 | 788 | 10 | | | |

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



| Mol | Chain | Residues | Atoms | | | ZeroOcc | AltConf |
|-----|-------|----------|-------|---|---|---------|---------|
| 2 | D | 1 | Total | C | O | 0 | 0 |
| | | | 6 | 3 | 3 | | |

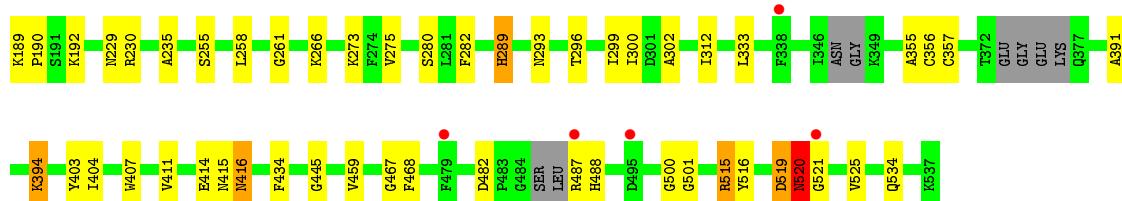
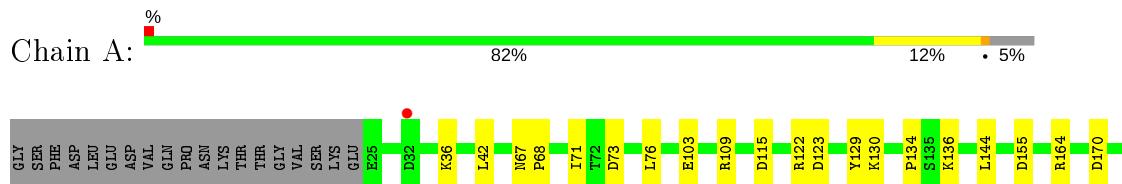
- Molecule 3 is water.

| Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|------------------|---------|---------|
| 3 | A | 79 | Total O 79 79 | 0 | 0 |
| 3 | B | 91 | Total O 91 91 | 0 | 0 |
| 3 | C | 48 | Total O 48 48 | 0 | 0 |
| 3 | D | 80 | Total O 80 80 | 0 | 0 |

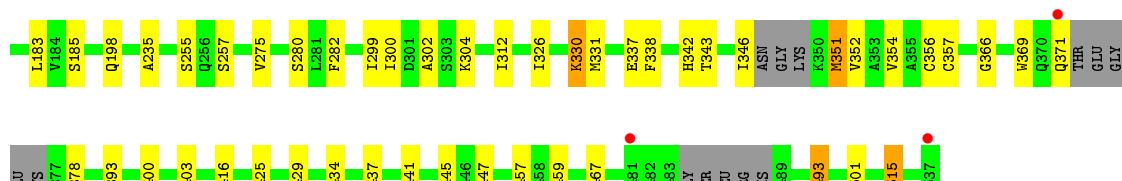
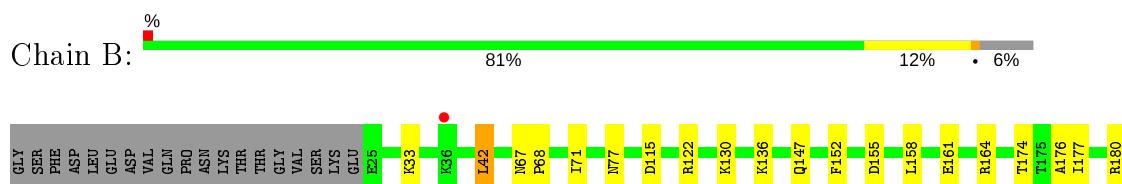
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

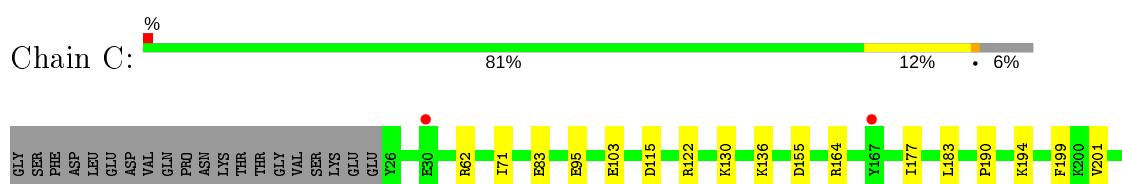
- Molecule 1: TbpB

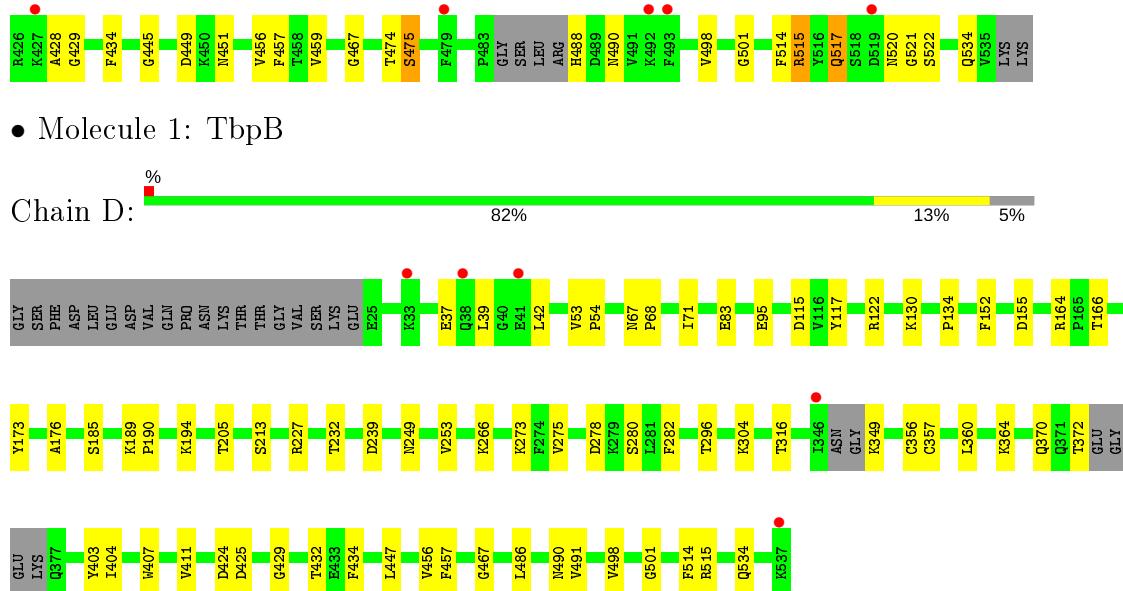


- Molecule 1: TbpB



- Molecule 1: TbpB





4 Data and refinement statistics i

| Property | Value | Source |
|---|--|------------------|
| Space group | P 1 21 1 | Depositor |
| Cell constants a, b, c, α , β , γ | 81.03 Å 135.70 Å 103.96 Å 90.00° 100.05° 90.00° | Depositor |
| Resolution (Å) | 48.39 – 2.64 48.39 – 2.64 | Depositor EDS |
| % Data completeness (in resolution range) | 99.5 (48.39-2.64) 74.3 (48.39-2.64) | Depositor EDS |
| R_{merge} | (Not available) | Depositor |
| R_{sym} | (Not available) | Depositor |
| $< I/\sigma(I) >$ ¹ | 2.07 (at 2.65 Å) | Xtriage |
| Refinement program | PHENIX 1.8.1_1168 | Depositor |
| R , R_{free} | 0.214 , 0.266 0.218 , 0.272 | Depositor DCC |
| R_{free} test set | 2393 reflections (4.96%) | wwPDB-VP |
| Wilson B-factor (Å ²) | 51.6 | Xtriage |
| Anisotropy | 0.272 | Xtriage |
| Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²) | 0.29 , 23.9 | EDS |
| L-test for twinning ² | $< L > = 0.50$, $< L^2 > = 0.33$ | Xtriage |
| Estimated twinning fraction | No twinning to report. | Xtriage |
| F_o, F_c correlation | 0.94 | EDS |
| Total number of atoms | 15892 | wwPDB-VP |
| Average B, all atoms (Å ²) | 34.0 | 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 37.79 % 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 4.0646e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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.

Continued from previous page...

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | A | 289 | HIS |
| 1 | A | 394 | LYS |
| 1 | A | 407 | TRP |
| 1 | A | 416 | ASN |
| 1 | A | 482 | ASP |
| 1 | A | 515 | ARG |
| 1 | A | 516 | TYR |
| 1 | A | 519 | ASP |
| 1 | A | 520 | ASN |
| 1 | B | 42 | LEU |
| 1 | B | 304 | LYS |
| 1 | B | 330 | LYS |
| 1 | B | 346 | ILE |
| 1 | B | 351 | MET |
| 1 | B | 493 | PHE |
| 1 | B | 515 | ARG |
| 1 | C | 62 | ARG |
| 1 | C | 83 | GLU |
| 1 | C | 95 | GLU |
| 1 | C | 103 | GLU |
| 1 | C | 304 | LYS |
| 1 | C | 346 | ILE |
| 1 | C | 394 | LYS |
| 1 | C | 398 | ASP |
| 1 | C | 488 | HIS |
| 1 | C | 490 | ASN |
| 1 | C | 515 | ARG |
| 1 | C | 517 | GLN |
| 1 | D | 37 | GLU |
| 1 | D | 39 | LEU |
| 1 | D | 42 | LEU |
| 1 | D | 213 | SER |
| 1 | D | 304 | LYS |
| 1 | D | 372 | THR |
| 1 | D | 407 | TRP |
| 1 | D | 486 | LEU |
| 1 | D | 490 | ASN |
| 1 | D | 515 | ARG |

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | C | 415 | ASN |
| 1 | D | 400 | 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 carbohydrates in this entry.

5.6 Ligand geometry [\(i\)](#)

1 ligand is 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 |
| 2 | GOL | D | 601 | - | 5,5,5 | 0.34 | 0 | 5,5,5 | 0.23 | 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 |
|-----|------|-------|-----|------|---------|----------|-------|
| 2 | GOL | D | 601 | - | - | 2/4/4/4 | - |

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

| Mol | Chain | Res | Type | Atoms |
|-----|-------|-----|------|-------------|
| 2 | D | 601 | GOL | O1-C1-C2-C3 |
| 2 | D | 601 | GOL | O1-C1-C2-O2 |

There are no ring outliers.

1 monomer is involved in 1 short contact:

| Mol | Chain | Res | Type | Clashes | Symm-Clashes |
|-----|-------|-----|------|---------|--------------|
| 2 | D | 601 | GOL | 1 | 0 |

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.

6.2 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. 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 | B-factors(Å ²) | Q<0.9 |
|-----|------|-------|-----|-------|------|------|----------------------------|-------|
| 2 | GOL | D | 601 | 6/6 | 0.91 | 0.18 | 37,41,45,48 | 0 |

6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.