

Full wwPDB X-ray Structure Validation Report (i)

Jun 25, 2024 – 04:51 AM EDT

PDB ID : 6POO

Title: Novel structure of the N-terminal helical domain of BibA, a group B strepto-

coccus immunogenic bacterial adhesin

Authors : Manne, K.; Narayana, S.V.

Deposited on : 2019-07-04

Resolution : 3.03 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 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:

 $Mol Probity \quad : \quad 4.02b\text{--}467$

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.37.1

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

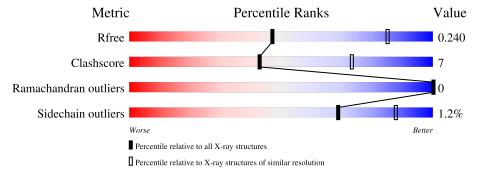
Validation Pipeline (wwPDB-VP) : 2.37.1

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 3.03 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	2752 (3.08-3.00)
Clashscore	141614	3096 (3.08-3.00)
Ramachandran outliers	138981	2986 (3.08-3.00)
Sidechain outliers	138945	2988 (3.08-3.00)

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%

16%
•



2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2158 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.

 \bullet Molecule 1 is a protein called BibA.

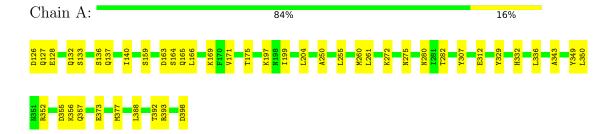
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	273	Total	С	N	О	Se	0	0	0
1	A	213	2158	1341	382	431	4	0	0	U



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. 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. 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: BibA





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 2 21	Depositor
Cell constants	26.42Å 68.81Å 199.38Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.84 - 3.03	Depositor
rtesolution (A)	68.81 - 1.80	EDS
% Data completeness	99.9 (49.84-3.03)	Depositor
(in resolution range)	63.5 (68.81-1.80)	EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	-0.05 (at 1.81Å)	Xtriage
Refinement program	PHENIX (1.15.2-3472)	Depositor
D D.	0.192 , 0.239	Depositor
R, R_{free}	0.192 , 0.240	DCC
R_{free} test set	1980 reflections (6.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 57.1	EDS
L-test for twinning ²	$ < L > = 0.46, < L^2> = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2158	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.57% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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).

Mal	Mol Chain Bond lengths		Bond	angles	
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.54	0/2172	0.61	0/2900

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 (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	2158	0	2228	30	1
All	All	2158	0	2228	30	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 7.

All (30) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:A:127:GLN:OE1	1:A:127:GLN:N	2.20	0.73
1:A:352:ARG:HD3	1:A:377:MSE:HE2	1.74	0.68
1:A:163:ASP:OD2	1:A:165:GLN:N	2.27	0.54
1:A:163:ASP:OD2	1:A:164:SER:N	2.41	0.53
1:A:352:ARG:NH1	1:A:377:MSE:HE1	2.24	0.53

Continued on next page...



Continued from previous page...

A. 1		Interatomic	Clash	
Atom-1	Atom-2	${\rm distance}\ ({\rm \AA})$	overlap (Å)	
1:A:137:GLN:OE1	1:A:159:SER:HB3	2.09	0.52	
1:A:329:TYR:CZ	1:A:343:ALA:HB2	2.47	0.50	
1:A:126:ASP:N	1:A:128:GLU:OE2	2.44	0.50	
1:A:165:GLN:O	1:A:169:LYS:HD2	2.12	0.50	
1:A:272:LYS:NZ	1:A:275:ASN:HD22	2.09	0.49	
1:A:352:ARG:HH11	1:A:377:MSE:HE1	1.76	0.49	
1:A:127:GLN:NE2	1:A:128:GLU:OE1	2.45	0.49	
1:A:133:SER:O	1:A:137:GLN:HG3	2.13	0.49	
1:A:250:ALA:HA	1:A:260:MSE:SE	2.62	0.49	
1:A:171:VAL:O	1:A:175:THR:HG23	2.14	0.48	
1:A:136:SER:O	1:A:140:ILE:HG12	2.14	0.47	
1:A:332:HIS:O	1:A:336:LEU:HD13	2.13	0.47	
1:A:349:TYR:CE1	1:A:377:MSE:HE3	2.51	0.45	
1:A:355:ASP:HB3	1:A:357:GLN:HG3	1.98	0.45	
1:A:197:LYS:HD2	1:A:197:LYS:HA	1.49	0.44	
1:A:388:LEU:O	1:A:392:THR:HG23	2.17	0.44	
1:A:352:ARG:O	1:A:356:LYS:N	2.48	0.43	
1:A:350:LEU:HD23	1:A:350:LEU:HA	1.74	0.43	
1:A:329:TYR:OH	1:A:343:ALA:HB2	2.19	0.43	
1:A:312:GLU:O	1:A:312:GLU:HG3	2.18	0.42	
1:A:280:ASN:OD1	1:A:282:THR:N	2.53	0.42	
1:A:128:GLU:O	1:A:132:GLN:N	2.38	0.42	
1:A:373:GLU:O	1:A:377:MSE:HG2	2.20	0.42	
1:A:199:ILE:HG12	1:A:204:LEU:HB2	2.01	0.41	
1:A:166:LEU:HA	1:A:169:LYS:HD3	2.03	0.41	

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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:255:LEU:O	1:A:393:ARG:NH2[3_577]	2.12	0.08

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	Perce	ntiles
1	A	271/273 (99%)	262 (97%)	9 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	240/236 (102%)	237 (99%)	3 (1%)	69 88

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	261	LEU
1	A	307	TYR
1	A	398	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 (i)

6.1 Protein, DNA and RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

