



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2018 – 12:10 pm GMT

PDB ID : 3Q7N
Title : The crystal structure of BamB from the BAM complex in spacegroup P212121
Authors : Noinaj, N.; Fairman, J.W.; Buchanan, S.K.
Deposited on : 2011-01-05
Resolution : 1.77 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : trunk31020
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk31020

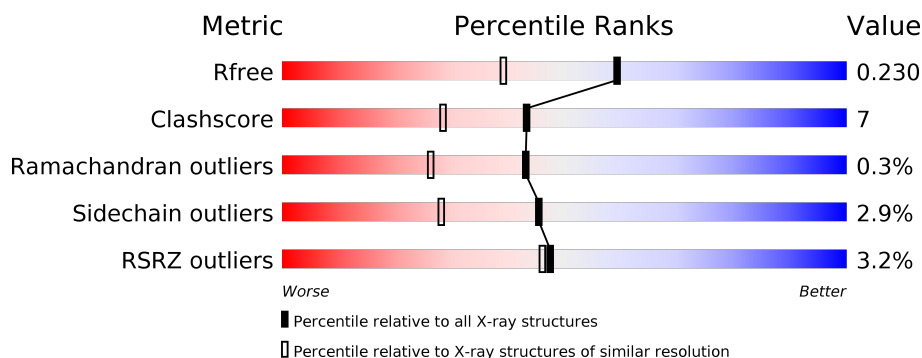
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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}	111664	7886 (1.80-1.76)
Clashscore	122126	8858 (1.80-1.76)
Ramachandran outliers	120053	8764 (1.80-1.76)
Sidechain outliers	120020	8763 (1.80-1.76)
RSRZ outliers	108989	7741 (1.80-1.76)

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. 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 $\leq 5\%$. 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	376	<div> <div>3%</div> <div> <div></div> <div>81%</div> <div>13%</div> <div>6%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2999 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 Lipoprotein yfgL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	355	Total	C	N	O	Se	0	0	0
			2639	1661	452	520	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	GLY	-	EXPRESSION TAG	UNP P77774
A	18	ALA	-	EXPRESSION TAG	UNP P77774
A	19	MSE	-	EXPRESSION TAG	UNP P77774
A	20	GLY	-	EXPRESSION TAG	UNP P77774

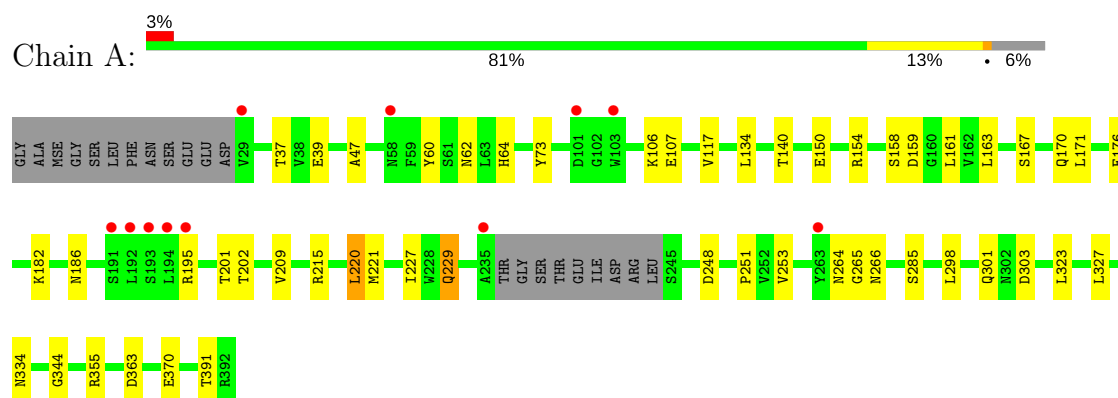
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	360	Total	O	0	0
			360	360		

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 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: Lipoprotein yfgL



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.00Å 69.99Å 117.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.62 – 1.77 30.62 – 1.77	Depositor EDS
% Data completeness (in resolution range)	84.8 (30.62-1.77) 84.5 (30.62-1.77)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 1.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_572)	Depositor
R, R_{free}	0.197 , 0.232 0.195 , 0.230	Depositor DCC
R_{free} test set	1702 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.506	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2999	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.59% 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, 0.333 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	0.36	0/2682	0.54	0/3651

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	2639	0	2560	38	0
2	A	360	0	0	8	1
All	All	2999	0	2560	38	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 (38) 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:140:THR:HA	2:A:663:HOH:O	1.88	0.73
1:A:62:ASN:O	1:A:64:HIS:HD2	1.73	0.71
1:A:220:LEU:HD22	1:A:227:ILE:HD11	1.74	0.70
1:A:201:THR:HG21	1:A:253:VAL:HG23	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:ARG:HD3	1:A:229:GLN:HE21	1.59	0.66
1:A:158:SER:HB3	1:A:221:MSE:HE3	1.76	0.66
1:A:215:ARG:HH11	1:A:229:GLN:NE2	1.96	0.63
1:A:170:GLN:HE21	1:A:186:ASN:HD22	1.47	0.61
1:A:171:LEU:HD22	1:A:209:VAL:HG21	1.84	0.60
1:A:154:ARG:HD2	2:A:532:HOH:O	2.03	0.59
1:A:163:LEU:HD11	1:A:221:MSE:HE1	1.87	0.57
1:A:201:THR:OG1	1:A:251:PRO:O	2.24	0.55
1:A:170:GLN:HE21	1:A:186:ASN:ND2	2.05	0.54
1:A:60:TYR:CD2	1:A:195:ARG:HB3	2.44	0.53
1:A:301:GLN:HG2	2:A:574:HOH:O	2.08	0.53
1:A:220:LEU:HD13	1:A:227:ILE:CG1	2.39	0.53
1:A:355:ARG:NH1	2:A:685:HOH:O	2.42	0.53
1:A:303:ASP:HB3	1:A:327:LEU:HG	1.90	0.52
1:A:47:ALA:HB2	1:A:391:THR:CG2	2.40	0.51
1:A:334:ASN:ND2	2:A:714:HOH:O	2.43	0.50
1:A:344:GLY:HA3	1:A:363:ASP:O	2.11	0.50
1:A:39:GLU:HG3	1:A:39:GLU:O	2.13	0.49
1:A:176:GLU:HG3	2:A:707:HOH:O	2.13	0.48
1:A:248:ASP:HB3	2:A:485:HOH:O	2.13	0.48
1:A:158:SER:HB3	1:A:221:MSE:CE	2.43	0.47
1:A:159:ASP:HA	2:A:690:HOH:O	2.16	0.46
1:A:220:LEU:CD2	1:A:227:ILE:HD11	2.44	0.46
1:A:220:LEU:HD13	1:A:227:ILE:HG13	1.98	0.45
1:A:150:GLU:HB2	1:A:167:SER:OG	2.16	0.45
1:A:220:LEU:HD13	1:A:227:ILE:HD11	2.00	0.44
1:A:264:ASN:OD1	1:A:265:GLY:N	2.52	0.43
1:A:298:LEU:C	1:A:298:LEU:HD12	2.39	0.43
1:A:158:SER:CB	1:A:221:MSE:HE3	2.48	0.42
1:A:161:LEU:HD21	1:A:182:LYS:HD2	2.01	0.42
1:A:201:THR:HG22	1:A:202:THR:N	2.35	0.42
1:A:47:ALA:HB2	1:A:391:THR:HG23	2.02	0.42
1:A:154:ARG:NH2	1:A:370:GLU:OE2	2.38	0.41
1:A:73:TYR:CD2	1:A:117:VAL:HG21	2.56	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	Interatomic distance (Å)	Clash overlap (Å)
2:A:623:HOH:O	2:A:687:HOH:O[1_455]	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	Percentiles	
1	A	351/376 (93%)	334 (95%)	16 (5%)	1 (0%)	43	26

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	285	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	277/297 (93%)	269 (97%)	8 (3%)	45	28

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	THR
1	A	106	LYS
1	A	107	GLU
1	A	134	LEU
1	A	220	LEU
1	A	229	GLN
1	A	266	ASN
1	A	323	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	HIS
1	A	121	HIS
1	A	135	ASN
1	A	186	ASN
1	A	213	ASN
1	A	229	GLN
1	A	255	ASN
1	A	266	ASN
1	A	302	ASN
1	A	324	HIS
1	A	336	ASN
1	A	359	GLN

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](#)

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](#)

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	349/376 (92%)	-0.08	11 (3%) 47 46	9, 17, 44, 72	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	191	SER	6.3
1	A	192	LEU	5.7
1	A	263	TYR	5.2
1	A	195	ARG	4.0
1	A	103	TRP	3.7
1	A	194	LEU	3.0
1	A	235	ALA	2.9
1	A	101	ASP	2.8
1	A	29	VAL	2.7
1	A	58	ASN	2.6
1	A	193	SER	2.6

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.