

wwPDB X-ray Structure Validation Summary Report (i)

Jul 1, 2024 – 08:10 PM JST

PDB ID : 8YK2

Title: Blood group B alpha-1,3-galactosidase AgaBb from Bifidobacterium bifidum,

construct T7-tag 24-700

Authors : Kashima, T.; Akama, M.; Ashida, H.; Fushinobu, S.

Deposited on : 2024-03-04

Resolution : 1.96 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:

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

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)

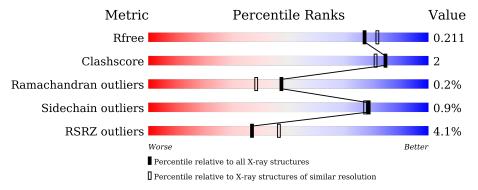
Validation Pipeline (wwPDB-VP) : 2.37.1

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 1.96 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}(\mathring{\rm A})) \end{array}$
R_{free}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	700	5% 87%		8%
1	В	700	87%	5%	7%



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 11006 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 Alpha-galactosidase.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	В	651	Total 4935	C 3082	N 838	O 1000	S 15	0	0	0
1	A	646	Total 4892	C 3057	N 831	O 989	S 15	0	0	0

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	9	MET	-	initiating methionine	UNP L8B3G2
В	10	ALA	-	expression tag	UNP L8B3G2
В	11	SER	=	expression tag	UNP L8B3G2
В	12	MET	-	expression tag	UNP L8B3G2
В	13	THR	-	expression tag	UNP L8B3G2
В	14	GLY	-	expression tag	UNP L8B3G2
В	15	GLY	-	expression tag	UNP L8B3G2
В	16	GLN	-	expression tag	UNP L8B3G2
В	17	GLN	-	expression tag	UNP L8B3G2
В	18	MET	-	expression tag	UNP L8B3G2
В	19	GLY	-	expression tag	UNP L8B3G2
В	20	ARG	-	expression tag	UNP L8B3G2
В	21	ASP	-	expression tag	UNP L8B3G2
В	22	PRO	-	expression tag	UNP L8B3G2
В	23	ASN	-	expression tag	UNP L8B3G2
В	701	LEU	-	expression tag	UNP L8B3G2
В	702	GLU	-	expression tag	UNP L8B3G2
В	703	HIS	-	expression tag	UNP L8B3G2
В	704	HIS	-	expression tag	UNP L8B3G2
В	705	HIS	-	expression tag	UNP L8B3G2
В	706	HIS	-	expression tag	UNP L8B3G2
В	707	HIS	-	expression tag	UNP L8B3G2
В	708	HIS	-	expression tag	UNP L8B3G2
A	9	MET	-	initiating methionine	UNP L8B3G2
A	10	ALA	-	expression tag	UNP L8B3G2

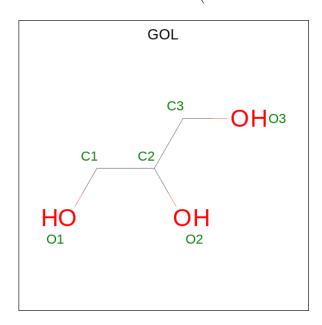
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Chain	Residue	Modelled	Actual	Comment	Reference
A	11	SER	-	expression tag	UNP L8B3G2
A	12	MET	-	expression tag	UNP L8B3G2
A	13	THR	-	expression tag	UNP L8B3G2
A	14	GLY	-	expression tag	UNP L8B3G2
A	15	GLY	-	expression tag	UNP L8B3G2
A	16	GLN	-	expression tag	UNP L8B3G2
A	17	GLN	-	expression tag	UNP L8B3G2
A	18	MET	-	expression tag	UNP L8B3G2
A	19	GLY	-	expression tag	UNP L8B3G2
A	20	ARG	-	expression tag	UNP L8B3G2
A	21	ASP	-	expression tag	UNP L8B3G2
A	22	PRO	-	expression tag	UNP L8B3G2
A	23	ASN	-	expression tag	UNP L8B3G2
A	701	LEU	-	expression tag	UNP L8B3G2
A	702	GLU	-	expression tag	UNP L8B3G2
A	703	HIS	-	expression tag	UNP L8B3G2
A	704	HIS	-	expression tag	UNP L8B3G2
A	705	HIS	-	expression tag	UNP L8B3G2
A	706	HIS	-	expression tag	UNP L8B3G2
A	707	HIS	-	expression tag	UNP L8B3G2
A	708	HIS	-	expression tag	UNP L8B3G2

 \bullet Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $\mathrm{C_3H_8O_3}).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	1	Total C O 6 3 3	0	0

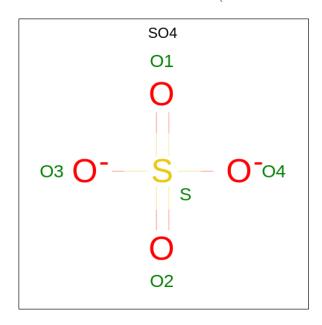
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	1	Total C O 6 3 3	0	0
2	В	1	Total C O 6 3 3	0	0

 \bullet Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 5	O 4	S 1	0	0

• Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Na 1 1	0	0

• Molecule 5 is water.

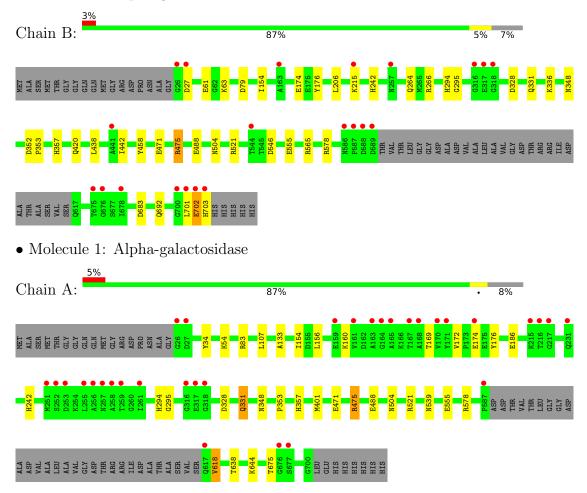
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	В	584	Total O 584 584	0	0
5	A	571	Total O 571 571	0	0



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: Alpha-galactosidase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	70.56Å 125.06Å 196.79Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.80 - 1.96	Depositor
rtesolution (A)	46.80 - 1.96	EDS
% Data completeness	100.0 (46.80-1.96)	Depositor
(in resolution range)	100.0 (46.80-1.96)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.16 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.21_5207	Depositor
D.D.	0.180 , 0.212	Depositor
R, R_{free}	0.179 , 0.211	DCC
R_{free} test set	6499 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	25.4	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.36, 40.9	EDS
L-test for twinning ²	$ < L > = 0.49, < L^2> = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11006	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.73% 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)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, NA, SO4

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		
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.35	0/4985	0.58	0/6777	
1	В	0.36	0/5029	0.59	0/6837	
All	All	0.35	0/10014	0.59	0/13614	

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	4892	0	4740	18	0
1	В	4935	0	4772	26	0
2	В	18	0	24	3	0
3	A	5	0	0	0	0
4	A	1	0	0	0	0
5	A	571	0	0	2	1
5	В	584	0	0	3	1
All	All	11006	0	9536	44	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 2.



The worst 5 of 44 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} & ext{Interatomic} \ & ext{distance} \ & ext{(Å)} \end{aligned}$	$egin{aligned} \operatorname{Clash} \ \operatorname{overlap}\ (\begin{subarray}{c} \begin{subarray}{c} \begi$
1:B:565:ARG:HH21	2:B:902:GOL:H31	1.49	0.77
1:B:692:GLN:NE2	5:B:1001:HOH:O	2.20	0.73
1:B:174:GLU:HG3	1:B:242:HIS:CE1	2.24	0.72
1:A:331:GLN:OE1	1:A:357:HIS:NE2	2.21	0.69
1:B:331:GLN:OE1	1:B:357:HIS:NE2	2.30	0.62

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)
5:B:1433:HOH:O	5:A:1778:HOH:O[3_655]	2.02	0.18

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	642/700 (92%)	623 (97%)	18 (3%)	1 (0%)	47	38
1	В	647/700 (92%)	625 (97%)	20 (3%)	2 (0%)	41	30
All	All	1289/1400 (92%)	1248 (97%)	38 (3%)	3 (0%)	47	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	701	LEU
1	В	504	ASN
1	A	504	ASN



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	Rotameric Outliers		Percentiles		
1	A	537/577 (93%)	532 (99%)	5 (1%)	78	77		
1	В	542/577 (94%)	537 (99%)	5 (1%)	78	77		
All	All	1079/1154 (94%)	1069 (99%)	10 (1%)	78	77		

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	331	GLN
1	A	475	ARG
1	A	618	VAL
1	В	546	ASP
1	В	702	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	420	GLN
1	A	350	HIS
1	В	651	ASN
1	В	348	ASN
1	A	348	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)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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	Trme	Chain	Res	Link	В	ond leng	gths	В	ond ang	gles
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	В	902	-	5,5,5	0.29	0	5,5,5	0.35	0
2	GOL	В	903	-	5,5,5	0.40	0	5,5,5	0.51	0
2	GOL	В	901	-	5,5,5	0.42	0	5,5,5	0.67	0
3	SO4	A	1301	-	4,4,4	0.59	0	6,6,6	0.07	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	В	902	_	-	0/4/4/4	-
2	GOL	В	903	-	-	2/4/4/4	_
2	GOL	В	901	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	901	GOL	O1-C1-C2-C3
2	В	903	GOL	C1-C2-C3-O3
2	В	901	GOL	O1-C1-C2-O2
2	В	903	GOL	O2-C2-C3-O3



There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	902	GOL	3	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 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, 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	$\langle { m RSRZ} \rangle$	# RSRZ > 2		$OWAB(A^2)$	Q < 0.9
1	A	646/700 (92%)	0.30	32 (4%) 2	28 39	15, 25, 49, 83	0
1	В	651/700 (93%)	0.08	21 (3%) 4	47 57	13, 24, 41, 95	0
All	All	1297/1400 (92%)	0.19	53 (4%) 3	37 46	13, 24, 46, 95	0

The worst 5 of 53 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	701	LEU	6.2
1	В	589	ASP	5.6
1	A	587	PRO	5.6
1	В	587	PRO	5.6
1	В	317	GLU	4.7

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 monosaccharides 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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	GOL	В	902	6/6	0.75	0.19	35,41,46,48	0
2	GOL	В	903	6/6	0.87	0.23	48,54,56,58	0
2	GOL	В	901	6/6	0.93	0.19	22,23,26,32	0
3	SO4	A	1301	5/5	0.96	0.18	47,52,56,59	0
4	NA	A	1302	1/1	0.96	0.34	44,44,44,44	0

6.5 Other polymers (i)

There are no such residues in this entry.

