



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:28 AM UTC

PDB ID : 8YK2 / pdb_00008yk2
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 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

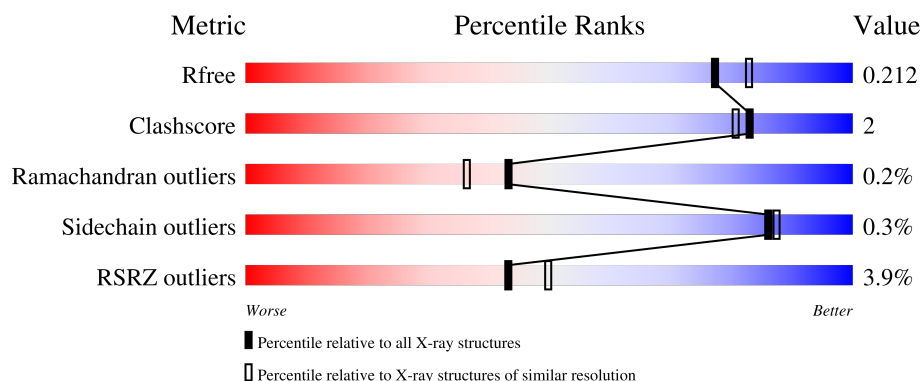
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

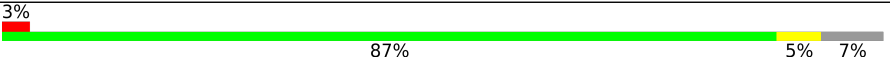
The reported resolution of this entry is 1.96 Å.

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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (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 $\leq 5\%$. 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	700	
1	B	700	

2 Entry composition

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	B	651	Total	C	N	O	S	0	0	0
			4935	3082	838	1000	15			
1	A	646	Total	C	N	O	S	0	0	0
			4892	3057	831	989	15			

There are 46 discrepancies between the modelled and reference sequences:

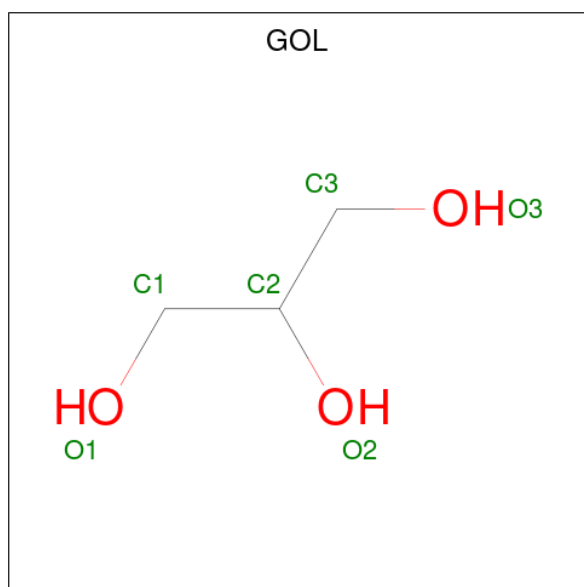
Chain	Residue	Modelled	Actual	Comment	Reference
B	9	MET	-	initiating methionine	UNP L8B3G2
B	10	ALA	-	expression tag	UNP L8B3G2
B	11	SER	-	expression tag	UNP L8B3G2
B	12	MET	-	expression tag	UNP L8B3G2
B	13	THR	-	expression tag	UNP L8B3G2
B	14	GLY	-	expression tag	UNP L8B3G2
B	15	GLY	-	expression tag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

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

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

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

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

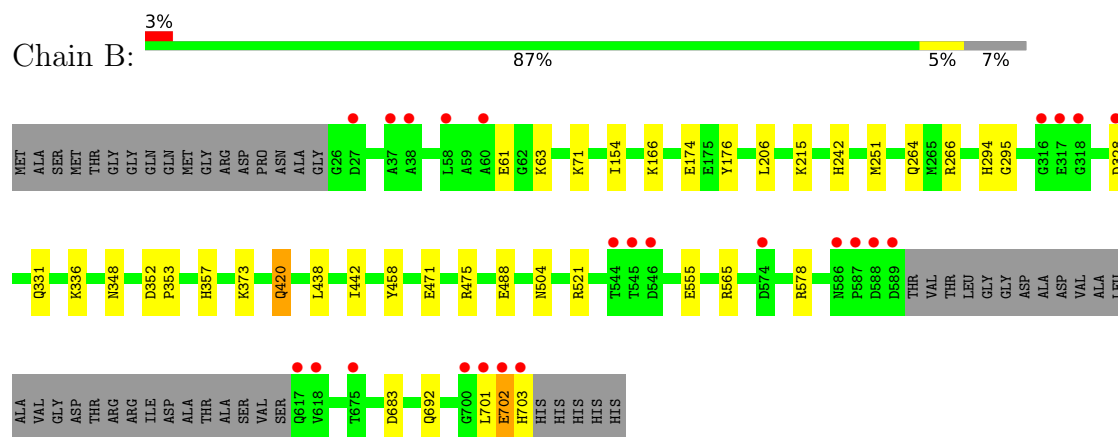
- Molecule 5 is water.

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

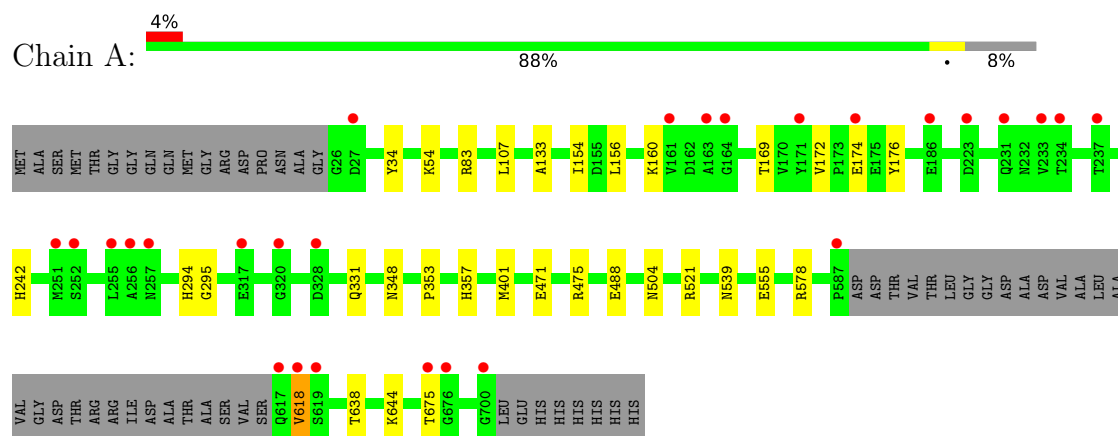
3 Residue-property plots [i](#)

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



- Molecule 1: Alpha-galactosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.56Å 125.06Å 196.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.80 – 1.96 46.80 – 1.96	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.80-1.96) 100.0 (46.80-1.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.21_5207	Depositor
R, R_{free}	0.180 , 0.212 0.180 , 0.212	Depositor 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 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	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.*

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

5.1 Standard geometry

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	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.27	0/4985	0.49	0/6777
1	B	0.28	0/5029	0.51	0/6837
All	All	0.27	0/10014	0.50	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

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	B	4935	0	4772	29	0
2	B	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	B	584	0	0	5	1
All	All	11006	0	9536	47	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.

All (47) 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: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
1:A:174:GLU:HG3	1:A:242:HIS:CE1	2.36	0.60
1:A:160:LYS:HG2	1:A:169:THR:HB	1.83	0.60
1:A:675:THR:HG22	5:A:1784:HOH:O	2.04	0.55
1:B:373:LYS:NZ	5:B:1002:HOH:O	2.35	0.55
1:B:206:LEU:HD23	1:B:266:ARG:HA	1.88	0.55
1:B:61:GLU:OE1	1:B:63:LYS:NZ	2.32	0.54
1:B:565:ARG:HH21	2:B:902:GOL:C3	2.20	0.53
1:B:264:GLN:HE22	1:B:348:ASN:ND2	2.06	0.52
1:B:353:PRO:HD2	1:B:475:ARG:O	2.12	0.50
1:B:264:GLN:HE22	1:B:348:ASN:HD21	1.61	0.49
1:A:353:PRO:HD2	1:A:475:ARG:O	2.11	0.49
1:B:215:LYS:HE2	5:B:1021:HOH:O	2.12	0.49
1:B:683:ASP:HA	1:B:703:HIS:CE1	2.49	0.48
1:B:521:ARG:HA	1:B:555:GLU:O	2.15	0.47
1:A:294:HIS:CG	1:A:295:GLY:N	2.83	0.47
1:B:336:LYS:HB3	1:B:458:TYR:HA	1.97	0.47
1:A:488:GLU:HA	1:A:521:ARG:O	2.14	0.47
1:A:83:ARG:HH11	1:A:83:ARG:HG2	1.80	0.46
1:B:348:ASN:HA	1:B:471:GLU:O	2.16	0.46
1:B:555:GLU:HA	1:B:578:ARG:O	2.17	0.45
1:A:348:ASN:HA	1:A:471:GLU:O	2.18	0.44
1:A:107:LEU:O	1:A:133:ALA:HA	2.17	0.44
1:A:401:MET:HG2	1:A:539:ASN:HB3	1.99	0.44
1:A:521:ARG:HA	1:A:555:GLU:O	2.18	0.44
1:B:328:ASP:OD1	1:B:331:GLN:HB2	2.19	0.43
1:B:328:ASP:HB2	1:B:352:ASP:O	2.18	0.43
1:B:71:LYS:NZ	5:B:1019:HOH:O	2.51	0.43
1:B:420:GLN:NE2	5:B:1014:HOH:O	2.49	0.43
1:B:294:HIS:CG	1:B:295:GLY:N	2.85	0.43
1:B:166:LYS:HA	1:B:251:MET:HE3	2.01	0.42
1:B:488:GLU:HA	1:B:521:ARG:O	2.20	0.42
1:A:156:LEU:HD22	1:A:172:VAL:HG22	2.02	0.42
1:B:438:LEU:HD12	1:B:442:ILE:CG2	2.51	0.41
1:B:565:ARG:NH2	2:B:902:GOL:H31	2.27	0.41
1:A:154:ILE:HA	1:A:176:TYR:CZ	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:ILE:HA	1:B:176:TYR:CZ	2.56	0.41
1:A:555:GLU:HA	1:A:578:ARG:O	2.21	0.41
1:A:638:THR:HG23	5:A:1619:HOH:O	2.20	0.41
1:A:34:TYR:OH	1:A:54:LYS:HG2	2.21	0.41
1:B:438:LEU:HD12	1:B:442:ILE:HG22	2.03	0.40
1:A:618:VAL:HB	1:A:644:LYS:HE2	2.02	0.40
1:B:702:GLU:H	1:B:702:GLU:HG3	1.66	0.40

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

5.3 Torsion angles

5.3.1 Protein backbone

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	642/700 (92%)	623 (97%)	18 (3%)	1 (0%)	43	36
1	B	647/700 (92%)	625 (97%)	20 (3%)	2 (0%)	36	28
All	All	1289/1400 (92%)	1248 (97%)	38 (3%)	3 (0%)	43	36

All (3) Ramachandran outliers are listed below:

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

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	537/577 (93%)	536 (100%)	1 (0%)	87	89
1	B	542/577 (94%)	540 (100%)	2 (0%)	84	85
All	All	1079/1154 (94%)	1076 (100%)	3 (0%)	86	87

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

Mol	Chain	Res	Type
1	B	420	GLN
1	B	702	GLU
1	A	618	VAL

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

Mol	Chain	Res	Type
1	B	64	ASN
1	B	177	ASN
1	B	232	ASN
1	B	257	ASN
1	B	343	ASN
1	B	348	ASN
1	B	489	ASN
1	A	267	GLN
1	A	348	ASN
1	A	350	HIS
1	A	420	GLN
1	A	651	ASN

5.3.3 RNA ⓘ

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 oligosaccharides 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	B	903	-	5,5,5	0.40	0	5,5,5	0.57	0
2	GOL	B	902	-	5,5,5	0.29	0	5,5,5	0.38	0
2	GOL	B	901	-	5,5,5	0.43	0	5,5,5	0.71	0
3	SO4	A	1301	-	4,4,4	0.66	0	6,6,6	0.09	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	B	903	-	-	2/4/4/4	-
2	GOL	B	902	-	-	0/4/4/4	-
2	GOL	B	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	B	901	GOL	O1-C1-C2-C3
2	B	903	GOL	C1-C2-C3-O3
2	B	901	GOL	O1-C1-C2-O2
2	B	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	B	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

6.1 Protein, DNA and RNA chains

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	646/700 (92%)	0.04	27 (4%)	40 47	15, 25, 49, 83	0
1	B	651/700 (93%)	0.00	24 (3%)	45 52	13, 24, 41, 95	0
All	All	1297/1400 (92%)	0.02	51 (3%)	43 50	13, 24, 46, 95	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	701	LEU	6.7
1	B	587	PRO	4.8
1	A	618	VAL	4.4
1	A	587	PRO	3.5
1	A	617	GLN	3.4
1	B	618	VAL	3.3
1	B	317	GLU	3.3
1	A	174	GLU	3.2
1	B	675	THR	3.2
1	A	251	MET	3.2
1	B	574	ASP	3.2
1	A	163	ALA	3.1
1	B	544	THR	3.1
1	A	675	THR	3.1
1	B	318	GLY	3.1
1	A	700	GLY	2.9
1	B	38	ALA	2.9
1	B	702	GLU	2.9
1	B	328	ASP	2.7
1	B	545	THR	2.7
1	A	161	VAL	2.7
1	A	255	LEU	2.6
1	A	237	THR	2.6
1	A	186	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	27	ASP	2.5
1	A	676	GLY	2.5
1	B	586	ASN	2.4
1	B	58	LEU	2.4
1	A	256	ALA	2.4
1	B	700	GLY	2.4
1	B	617	GLN	2.3
1	B	703	HIS	2.3
1	B	589	ASP	2.3
1	A	223	ASP	2.3
1	A	233	VAL	2.3
1	A	234	THR	2.3
1	A	257	ASN	2.3
1	A	619	SER	2.2
1	A	252	SER	2.2
1	B	546	ASP	2.2
1	A	171	TYR	2.1
1	A	320	GLY	2.1
1	B	60	ALA	2.1
1	B	316	GLY	2.1
1	B	37	ALA	2.1
1	B	588	ASP	2.1
1	A	27	ASP	2.1
1	A	328	ASP	2.1
1	A	164	GLY	2.0
1	A	231	GLN	2.0
1	A	317	GLU	2.0

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	GOL	B	902	6/6	0.73	0.16	35,41,46,48	0
2	GOL	B	903	6/6	0.73	0.24	48,54,56,58	0
2	GOL	B	901	6/6	0.90	0.11	22,23,26,32	0
4	NA	A	1302	1/1	0.91	0.38	44,44,44,44	0
3	SO4	A	1301	5/5	0.93	0.07	47,52,56,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.