



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:53 AM UTC

PDB ID : 8YK1 / pdb_00008yk1
Title : Blood group B alpha-1,3-galactosidase AgaBb from Bifidobacterium bifidum, construct 23-844
Authors : Kashima, T.; Ashida, H.; Fushinobu, S.
Deposited on : 2024-03-04
Resolution : 2.02 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
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

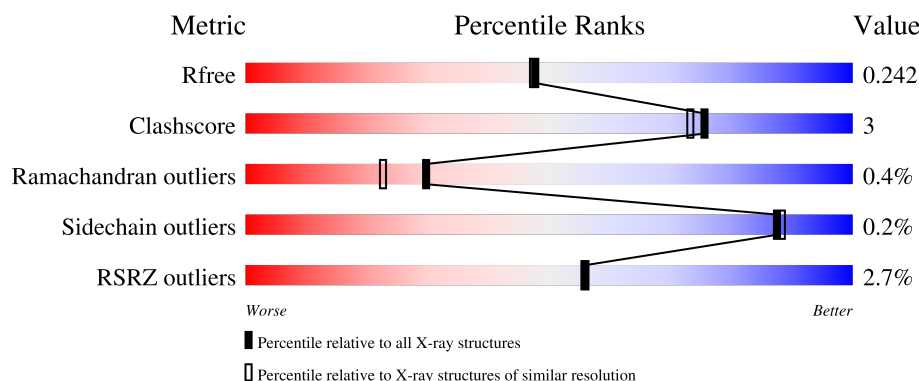
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

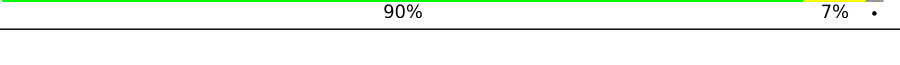
The reported resolution of this entry is 2.02 Å.

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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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 $\leq 5\%$. 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	829	
1	B	829	
1	C	829	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 19153 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	A	810	Total	C	N	O	S	0	0	0
			6072	3783	1040	1233	16			
1	B	810	Total	C	N	O	S	0	0	0
			6072	3783	1040	1233	16			
1	C	810	Total	C	N	O	S	0	0	0
			6072	3783	1040	1233	16			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	initiating methionine	UNP L8B3G2
A	845	LEU	-	expression tag	UNP L8B3G2
A	846	GLU	-	expression tag	UNP L8B3G2
A	847	HIS	-	expression tag	UNP L8B3G2
A	848	HIS	-	expression tag	UNP L8B3G2
A	849	HIS	-	expression tag	UNP L8B3G2
A	850	HIS	-	expression tag	UNP L8B3G2
A	851	HIS	-	expression tag	UNP L8B3G2
A	852	HIS	-	expression tag	UNP L8B3G2
B	24	MET	-	initiating methionine	UNP L8B3G2
B	845	LEU	-	expression tag	UNP L8B3G2
B	846	GLU	-	expression tag	UNP L8B3G2
B	847	HIS	-	expression tag	UNP L8B3G2
B	848	HIS	-	expression tag	UNP L8B3G2
B	849	HIS	-	expression tag	UNP L8B3G2
B	850	HIS	-	expression tag	UNP L8B3G2
B	851	HIS	-	expression tag	UNP L8B3G2
B	852	HIS	-	expression tag	UNP L8B3G2
C	24	MET	-	initiating methionine	UNP L8B3G2
C	845	LEU	-	expression tag	UNP L8B3G2
C	846	GLU	-	expression tag	UNP L8B3G2
C	847	HIS	-	expression tag	UNP L8B3G2
C	848	HIS	-	expression tag	UNP L8B3G2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	849	HIS	-	expression tag	UNP L8B3G2
C	850	HIS	-	expression tag	UNP L8B3G2
C	851	HIS	-	expression tag	UNP L8B3G2
C	852	HIS	-	expression tag	UNP L8B3G2

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Na 1 1	0	0
2	B	1	Total Na 1 1	0	0
2	C	1	Total Na 1 1	0	0

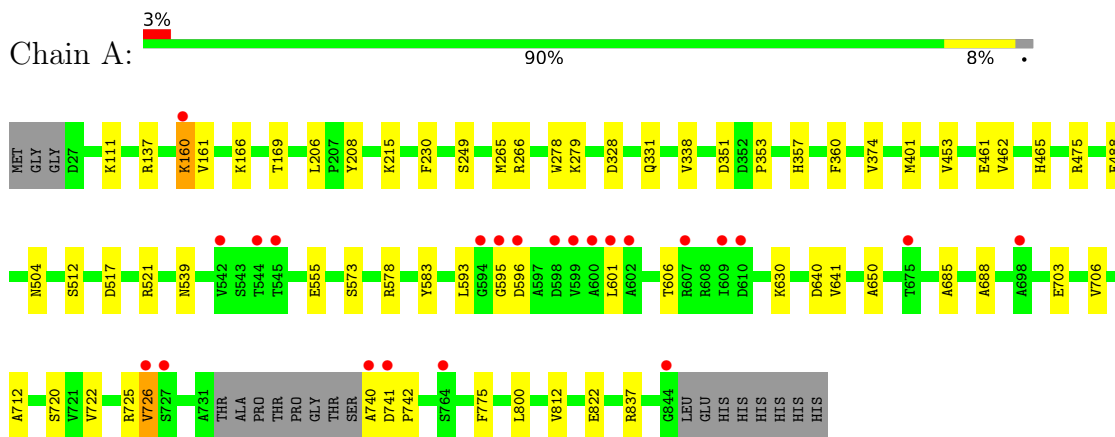
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	357	Total O 357 357	0	0
3	B	368	Total O 368 368	0	0
3	C	209	Total O 209 209	0	0

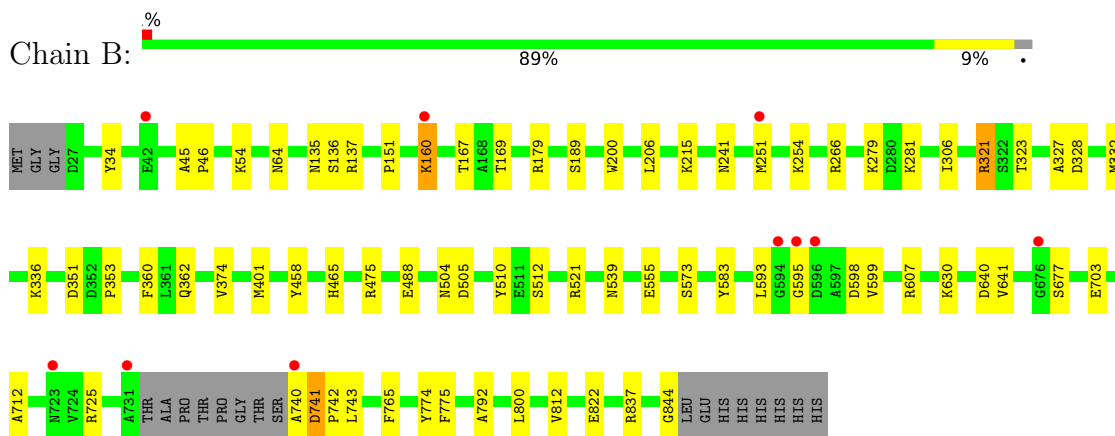
3 Residue-property plots [i](#)

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.

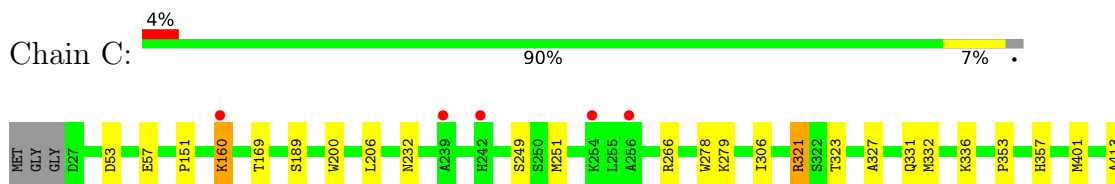
- Molecule 1: Alpha-galactosidase

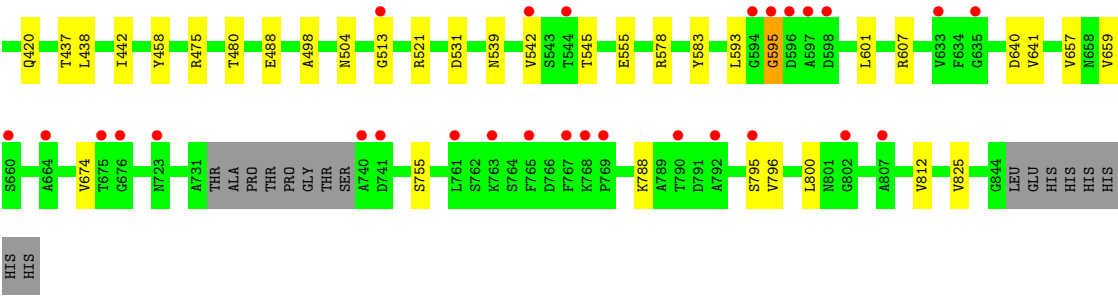


- Molecule 1: Alpha-galactosidase



- Molecule 1: Alpha-galactosidase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	127.69Å 68.77Å 291.31Å 90.00° 91.40° 90.00°	Depositor
Resolution (Å)	48.59 – 2.02 48.59 – 2.02	Depositor EDS
% Data completeness (in resolution range)	98.0 (48.59-2.02) 97.4 (48.59-2.02)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.03Å)	Xtriage
Refinement program	PHENIX 1.21_5207	Depositor
R, R_{free}	0.189 , 0.221 (Not available) , 0.242	Depositor DCC
R_{free} test set	8190 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.004 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	19153	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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:
NA

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	0.33	0/6179	0.53	1/8405 (0.0%)
1	B	0.32	1/6179 (0.0%)	0.51	0/8405
1	C	0.26	1/6179 (0.0%)	0.45	0/8405
All	All	0.30	2/18537 (0.0%)	0.50	1/25215 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	160	LYS	CE-NZ	-6.79	1.28	1.49
1	B	160	LYS	CE-NZ	-5.74	1.32	1.49

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	160	LYS	CD-CE-NZ	6.01	131.14	111.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	321	ARG	Sidechain
1	C	321	ARG	Sidechain

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	6072	0	5911	37	0
1	B	6072	0	5911	42	0
1	C	6072	0	5911	31	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	357	0	0	0	0
3	B	368	0	0	1	0
3	C	209	0	0	1	0
All	All	19153	0	17733	108	0

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 3.

All (108) 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:137:ARG:HG3	1:B:281:LYS:HB3	1.68	0.76
1:B:179:ARG:NH2	3:B:1001:HOH:O	2.23	0.71
1:A:703:GLU:HG2	1:A:725:ARG:HG2	1.71	0.71
1:A:331:GLN:OE1	1:A:357:HIS:NE2	2.23	0.68
1:A:800:LEU:HD22	1:A:812:VAL:HB	1.77	0.67
1:B:598:ASP:OD2	1:B:725:ARG:NH1	2.29	0.66
1:C:800:LEU:HD22	1:C:812:VAL:HB	1.78	0.65
1:C:796:VAL:HG22	1:C:825:VAL:HG22	1.79	0.64
1:A:215:LYS:HG2	1:A:775:PHE:HB2	1.80	0.62
1:B:703:GLU:HG3	1:B:725:ARG:HG2	1.83	0.61
1:B:241:ASN:ND2	1:B:844:GLY:O	2.33	0.61
1:A:160:LYS:HE3	1:A:169:THR:OG1	2.02	0.60
1:B:215:LYS:HG2	1:B:775:PHE:HB2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:MET:HG2	1:A:539:ASN:HB3	1.84	0.59
1:B:306:ILE:HD12	1:B:332:MET:SD	2.43	0.58
1:C:321:ARG:NH2	1:C:323:THR:O	2.36	0.58
1:C:206:LEU:HD23	1:C:266:ARG:HA	1.85	0.57
1:C:401:MET:HG2	1:C:539:ASN:HB3	1.87	0.57
1:B:160:LYS:HG3	1:B:169:THR:HB	1.87	0.57
1:A:353:PRO:HD2	1:A:475:ARG:O	2.06	0.56
1:A:230:PHE:HE2	1:A:265:MET:HE1	1.71	0.55
1:B:321:ARG:NH2	1:B:323:THR:O	2.40	0.55
1:B:593:LEU:O	1:B:595:GLY:N	2.38	0.54
1:C:601:LEU:HD21	1:C:607:ARG:HB2	1.89	0.54
1:B:353:PRO:HD2	1:B:475:ARG:O	2.08	0.54
1:B:206:LEU:HD23	1:B:266:ARG:HA	1.90	0.53
1:B:800:LEU:HD22	1:B:812:VAL:HB	1.90	0.53
1:C:151:PRO:HG2	1:C:321:ARG:NE	2.25	0.52
1:A:328:ASP:OD1	1:A:351:ASP:HB3	2.10	0.52
1:A:338:VAL:HG13	1:A:461:GLU:HG2	1.91	0.52
1:C:583:TYR:HA	1:C:640:ASP:HB2	1.91	0.52
1:A:601:LEU:HB2	1:A:726:VAL:HG12	1.91	0.52
1:B:599:VAL:HG13	1:B:607:ARG:HD3	1.91	0.51
1:B:521:ARG:HA	1:B:555:GLU:O	2.11	0.51
1:A:740:ALA:HB1	1:A:742:PRO:HD2	1.93	0.51
1:B:360:PHE:CE2	1:B:512:SER:HA	2.46	0.51
1:B:136:SER:C	1:B:137:ARG:HD2	2.35	0.51
1:B:254:LYS:HE2	1:B:254:LYS:H	1.76	0.51
1:C:755:SER:HB3	1:C:788:LYS:HB2	1.93	0.50
1:B:740:ALA:HB1	1:B:743:LEU:HB2	1.94	0.49
1:C:336:LYS:HB3	1:C:458:TYR:HA	1.94	0.49
1:C:795:SER:O	1:C:825:VAL:HA	2.12	0.49
1:A:360:PHE:CE2	1:A:512:SER:HA	2.48	0.48
1:B:488:GLU:HA	1:B:521:ARG:O	2.13	0.48
1:C:521:ARG:HA	1:C:555:GLU:O	2.13	0.48
1:B:251:MET:HE3	1:B:251:MET:HB2	1.68	0.48
1:A:583:TYR:HA	1:A:640:ASP:HB2	1.94	0.48
1:C:420:GLN:NE2	3:C:1003:HOH:O	2.41	0.48
1:C:251:MET:HE2	1:C:251:MET:HB3	1.69	0.48
1:C:353:PRO:HD2	1:C:475:ARG:O	2.13	0.48
1:B:583:TYR:HA	1:B:640:ASP:HB2	1.96	0.47
1:A:606:THR:OG1	1:B:64:ASN:ND2	2.47	0.47
1:C:641:VAL:CG1	1:C:674:VAL:HA	2.43	0.47
1:C:593:LEU:C	1:C:595:GLY:H	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:LYS:HA	1:A:137:ARG:O	2.15	0.47
1:B:741:ASP:OD1	1:B:742:PRO:HD3	2.15	0.46
1:B:160:LYS:HE3	1:B:169:THR:OG1	2.15	0.46
1:A:461:GLU:HG3	1:A:462:VAL:N	2.32	0.45
1:C:413:ALA:HB3	1:C:437:THR:HB	1.97	0.45
1:C:555:GLU:HA	1:C:578:ARG:O	2.16	0.45
1:C:488:GLU:HA	1:C:521:ARG:O	2.17	0.45
1:A:488:GLU:HA	1:A:521:ARG:O	2.17	0.45
1:B:822:GLU:HG2	1:B:837:ARG:HG2	1.99	0.45
1:B:151:PRO:HG2	1:B:321:ARG:NE	2.32	0.44
1:B:189:SER:HB3	1:B:200:TRP:CE2	2.52	0.44
1:C:160:LYS:HG3	1:C:169:THR:HB	2.00	0.44
1:A:706:VAL:HG22	1:A:722:VAL:HB	1.98	0.44
1:A:166:LYS:HE3	1:A:249:SER:O	2.17	0.44
1:C:657:VAL:HG12	1:C:659:VAL:HG13	1.98	0.44
1:C:53:ASP:O	1:C:57:GLU:HG3	2.18	0.44
1:A:641:VAL:HG21	1:A:712:ALA:HB1	1.98	0.44
1:B:328:ASP:OD1	1:B:351:ASP:HB3	2.18	0.44
1:C:278:TRP:CE2	1:C:279:LYS:HE2	2.52	0.44
1:C:189:SER:HB3	1:C:200:TRP:CE2	2.54	0.43
1:A:208:TYR:O	1:A:265:MET:HA	2.19	0.43
1:A:521:ARG:HA	1:A:555:GLU:O	2.17	0.43
1:A:465:HIS:HA	1:A:488:GLU:O	2.19	0.43
1:B:160:LYS:NZ	1:B:167:THR:OG1	2.51	0.43
1:B:641:VAL:HG21	1:B:712:ALA:HB1	2.00	0.43
1:B:465:HIS:HA	1:B:488:GLU:O	2.19	0.43
1:C:480:THR:O	1:C:513:GLY:HA3	2.19	0.42
1:A:573:SER:HA	1:A:630:LYS:O	2.18	0.42
1:C:232:ASN:HB3	1:C:249:SER:OG	2.20	0.42
1:C:331:GLN:OE1	1:C:357:HIS:NE2	2.39	0.42
1:A:685:ALA:HA	1:A:688:ALA:O	2.20	0.42
1:B:573:SER:HA	1:B:630:LYS:O	2.19	0.41
1:A:206:LEU:HB2	1:A:265:MET:HG3	2.01	0.41
1:B:362:GLN:O	1:B:374:VAL:HA	2.20	0.41
1:B:45:ALA:HB3	1:B:46:PRO:HD3	2.02	0.41
1:B:401:MET:HG2	1:B:539:ASN:HB3	2.02	0.41
1:A:206:LEU:HD23	1:A:266:ARG:HA	2.02	0.41
1:A:278:TRP:CE2	1:A:279:LYS:HE2	2.55	0.41
1:A:822:GLU:HG2	1:A:837:ARG:HG2	2.03	0.41
1:A:593:LEU:HG	1:A:720:SER:HB2	2.03	0.41
1:A:650:ALA:HB1	1:B:792:ALA:HB1	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ASN:OD1	1:B:279:LYS:HE3	2.20	0.41
1:B:765:PHE:CD1	1:B:774:TYR:HB3	2.55	0.41
1:A:555:GLU:HA	1:A:578:ARG:O	2.20	0.41
1:A:741:ASP:N	1:A:742:PRO:HD2	2.36	0.41
1:B:505:ASP:O	1:B:510:TYR:HA	2.21	0.41
1:C:306:ILE:HD12	1:C:332:MET:SD	2.61	0.40
1:A:374:VAL:HG11	1:A:453:VAL:HG11	2.02	0.40
1:C:498:ALA:HA	1:C:531:ASP:O	2.21	0.40
1:A:160:LYS:HG2	1:A:161:VAL:N	2.35	0.40
1:C:438:LEU:HD12	1:C:442:ILE:CG2	2.51	0.40
1:A:593:LEU:O	1:A:595:GLY:N	2.52	0.40
1:B:34:TYR:OH	1:B:54:LYS:HG2	2.21	0.40
1:B:336:LYS:HB3	1:B:458:TYR:HA	2.02	0.40

There are no symmetry-related clashes.

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	806/829 (97%)	777 (96%)	26 (3%)	3 (0%)	30	22
1	B	806/829 (97%)	772 (96%)	31 (4%)	3 (0%)	30	22
1	C	806/829 (97%)	772 (96%)	31 (4%)	3 (0%)	30	22
All	All	2418/2487 (97%)	2321 (96%)	88 (4%)	9 (0%)	30	22

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	596	ASP
1	B	677	SER
1	B	504	ASN
1	A	504	ASN

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Mol	Chain	Res	Type
1	A	517	ASP
1	C	327	ALA
1	C	504	ASN
1	C	595	GLY
1	B	327	ALA

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	658/674 (98%)	657 (100%)	1 (0%)	87	88
1	B	658/674 (98%)	657 (100%)	1 (0%)	87	88
1	C	658/674 (98%)	656 (100%)	2 (0%)	86	86
All	All	1974/2022 (98%)	1970 (100%)	4 (0%)	87	88

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	726	VAL
1	B	741	ASP
1	C	542	VAL
1	C	545	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	489	ASN
1	A	617	GLN
1	B	64	ASN
1	C	64	ASN
1	C	231	GLN
1	C	489	ASN
1	C	617	GLN
1	C	746	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	810/829 (97%)	-0.16	23 (2%) 55 54	12, 21, 50, 75	0
1	B	810/829 (97%)	-0.21	10 (1%) 76 76	12, 23, 38, 58	0
1	C	810/829 (97%)	0.24	33 (4%) 41 41	17, 31, 50, 68	0
All	All	2430/2487 (97%)	-0.04	66 (2%) 56 56	12, 24, 47, 75	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	740	ALA	6.8
1	A	740	ALA	5.1
1	B	676	GLY	4.5
1	A	594	GLY	4.1
1	C	740	ALA	3.9
1	C	542	VAL	3.9
1	A	542	VAL	3.8
1	C	595	GLY	3.5
1	A	160	LYS	3.5
1	B	160	LYS	3.4
1	C	763	LYS	3.3
1	C	160	LYS	3.3
1	A	595	GLY	3.3
1	C	675	THR	3.2
1	A	675	THR	3.1
1	C	768	LYS	3.1
1	A	602	ALA	3.0
1	A	599	VAL	3.0
1	B	596	ASP	3.0
1	C	795	SER	2.8
1	B	594	GLY	2.8
1	C	767	PHE	2.7
1	A	600	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	596	ASP	2.7
1	C	513	GLY	2.7
1	C	597	ALA	2.6
1	C	741	ASP	2.5
1	C	239	ALA	2.5
1	C	598	ASP	2.5
1	C	676	GLY	2.5
1	A	609	ILE	2.5
1	B	595	GLY	2.5
1	A	727	SER	2.4
1	B	731	ALA	2.4
1	C	792	ALA	2.4
1	B	723	ASN	2.4
1	A	610	ASP	2.3
1	A	741	ASP	2.3
1	C	256	ALA	2.3
1	A	607	ARG	2.3
1	C	254	LYS	2.3
1	C	596	ASP	2.3
1	B	251	MET	2.3
1	A	764	SER	2.2
1	A	545	THR	2.2
1	C	765	PHE	2.2
1	A	544	THR	2.2
1	C	807	ALA	2.2
1	C	544	THR	2.2
1	A	844	GLY	2.1
1	A	601	LEU	2.1
1	C	761	LEU	2.1
1	C	723	ASN	2.1
1	A	726	VAL	2.1
1	C	594	GLY	2.1
1	C	633	VAL	2.1
1	C	242	HIS	2.1
1	B	42	GLU	2.1
1	A	598	ASP	2.1
1	C	790	THR	2.1
1	C	635	GLY	2.1
1	C	664	ALA	2.1
1	C	660	SER	2.0
1	A	698	ALA	2.0
1	C	769	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	802	GLY	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(Å ²)	Q<0.9
2	NA	A	901	1/1	0.95	0.25	14,14,14,14	0
2	NA	B	901	1/1	0.98	0.29	3,3,3,3	0
2	NA	C	901	1/1	0.99	0.29	3,3,3,3	0

6.5 Other polymers [i](#)

There are no such residues in this entry.