Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2026 – 04:14 PM JST

PDB ID	:	9XKP / pdb_00009xkp
Title	:	Tepidibacillus decaturensis alpha-glucan 4(6)-alpha-glucosyltransferase E666Q in apo-form
Authors	:	Qin, T.; Saburi, W.; Takei, S.; Tagami, T.; Ota, T.; Kanai, K.; Ose, T.; Svensson, B.; Mori, H.
Deposited on	:	2025-11-06
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 (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 (i)) were used in the production of this report:

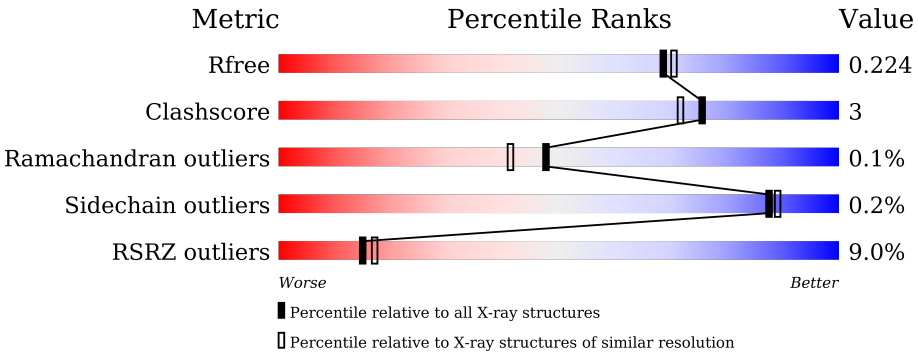
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

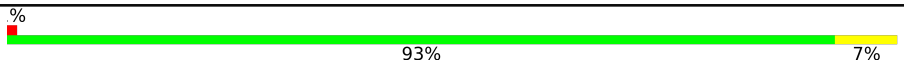
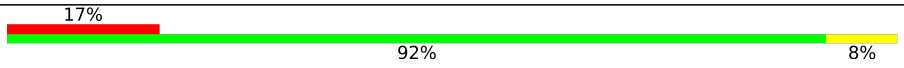
The reported resolution of this entry is 1.96 Å.

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 (#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 <=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	887	
1	B	887	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15143 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.

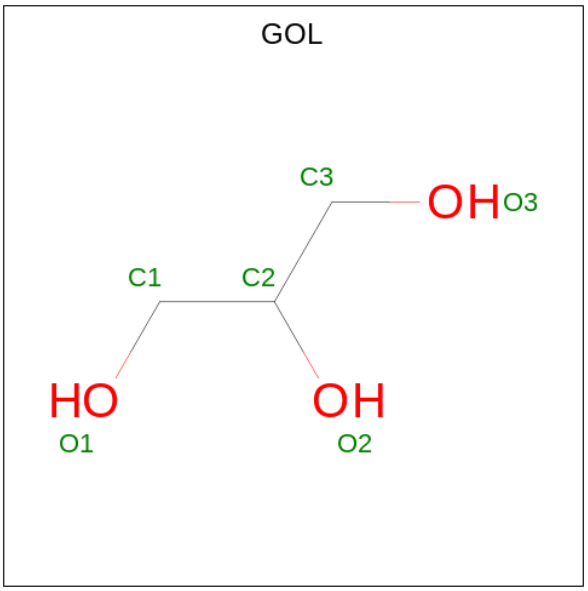
- Molecule 1 is a protein called GH15-like domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	887	Total	C	N	O	S	0	0	0
			7016	4414	1168	1418	16			
1	B	887	Total	C	N	O	S	0	0	0
			7016	4414	1168	1418	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	666	GLN	GLU	engineered mutation	UNP A0A135L1V7
B	666	GLN	GLU	engineered mutation	UNP A0A135L1V7

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



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

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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	5	Total Ca 5 5	0	0
3	B	5	Total Ca 5 5	0	0

- Molecule 4 is water.

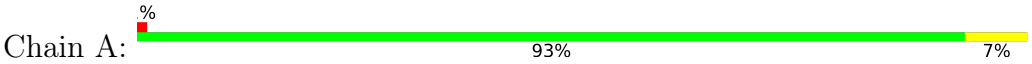
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	615	Total O 615 615	0	0
4	B	444	Total O 444 444	0	0

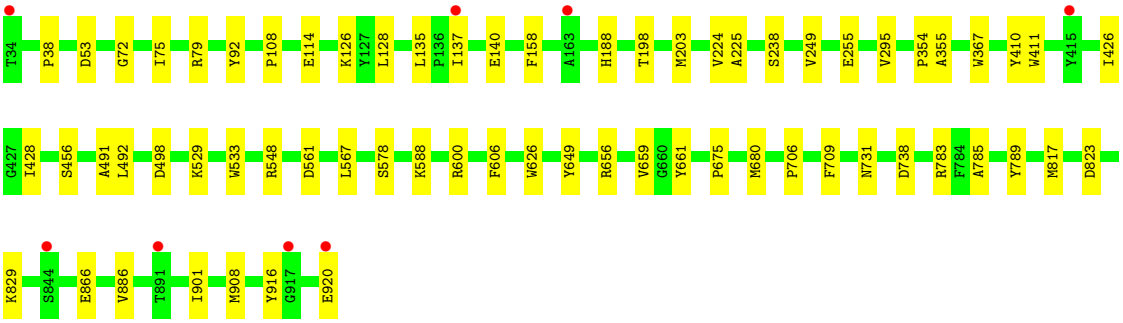
3 Residue-property plots (i)

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: GH15-like domain-containing protein

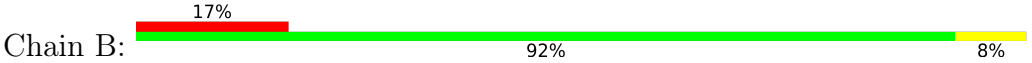
Chain A:

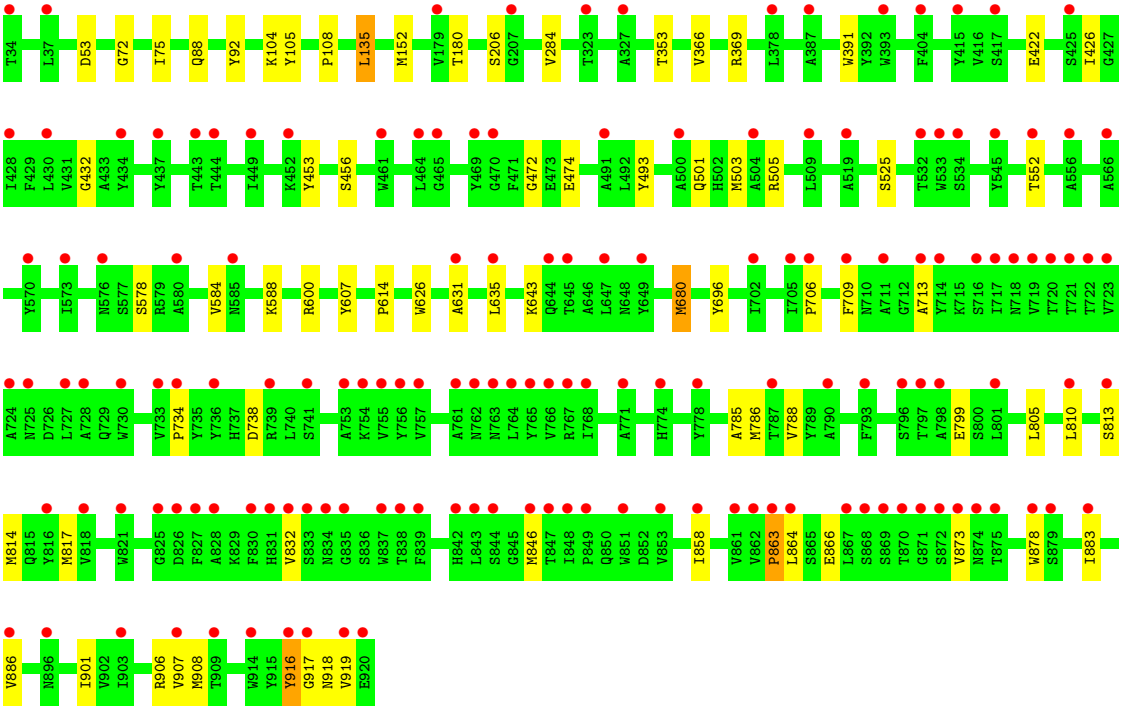




• Molecule 1: GH15-like domain-containing protein

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	112.92Å 112.92Å 391.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.80 – 1.96 48.80 – 1.96	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.80-1.96) 100.0 (48.80-1.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.64 (at 1.86Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.211 , 0.229 0.206 , 0.224	Depositor DCC
R_{free} test set	2000 reflections (0.93%)	wwPDB-VP
Wilson B-factor (Å^2)	27.5	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 38.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	15143	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.67% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

5 Model quality

5.1 Standard geometry

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

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	0.25	0/7200	0.49	2/9839 (0.0%)
1	B	0.28	1/7200 (0.0%)	0.51	5/9839 (0.1%)
All	All	0.26	1/14400 (0.0%)	0.50	7/19678 (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	A	0	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	863	PRO	CG-CD	-9.47	1.18	1.50

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	863	PRO	N-CD-CG	-16.25	78.83	103.20
1	B	863	PRO	CA-CB-CG	-13.39	79.06	104.50
1	B	680	MET	CA-C-N	5.72	131.99	121.70
1	B	680	MET	C-N-CA	5.72	131.99	121.70
1	B	863	PRO	N-CA-CB	-5.70	98.14	103.27
1	A	680	MET	CA-C-N	5.08	130.85	121.70
1	A	680	MET	C-N-CA	5.08	130.85	121.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	916	TYR	Peptide
1	B	916	TYR	Peptide

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	7016	0	6505	38	0
1	B	7016	0	6505	42	0
2	A	30	0	40	6	0
2	B	12	0	16	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	615	0	0	0	0
4	B	444	0	0	6	0
All	All	15143	0	13066	80	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 (80) 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:135:LEU:H	2:A:1001:GOL:H31	1.21	1.02
1:A:249:VAL:H	2:A:1002:GOL:H31	1.34	0.92
1:B:786:MET:HG2	1:B:858:ILE:HD13	1.70	0.71
1:B:738:ASP:HB2	1:B:901:ILE:HD12	1.76	0.67
1:B:72:GLY:HA3	1:B:108:PRO:HB2	1.80	0.64
1:B:706:PRO:HG2	1:B:709:PHE:CE2	2.34	0.62
1:A:731:ASN:O	1:A:920:GLU:HG2	1.99	0.62
1:B:426:ILE:HG23	1:B:456:SER:HB3	1.83	0.61
1:A:785:ALA:HB3	1:A:886:VAL:HB	1.83	0.59
1:A:72:GLY:HA3	1:A:108:PRO:HB2	1.85	0.59
1:A:188:HIS:CE1	1:A:203:MET:HG2	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:369:ARG:NH1	1:B:422:GLU:OE2	2.37	0.58
1:A:135:LEU:N	2:A:1001:GOL:H31	2.06	0.57
1:B:863:PRO:HB2	1:B:866:GLU:HG2	1.86	0.57
1:A:79:ARG:HD2	1:A:411:TRP:CH2	2.41	0.56
1:B:917:GLY:HA2	4:B:1476:HOH:O	2.06	0.55
1:B:785:ALA:HB3	1:B:886:VAL:HB	1.88	0.54
1:A:738:ASP:HB2	1:A:901:ILE:HD12	1.90	0.53
1:B:584:VAL:O	1:B:588:LYS:HG3	2.08	0.53
1:B:75:ILE:HB	1:B:92:TYR:HB3	1.92	0.52
1:A:656:ARG:O	1:A:675:PRO:HG3	2.09	0.51
1:B:680:MET:HB3	4:B:1270:HOH:O	2.10	0.51
1:B:643:LYS:NZ	4:B:1104:HOH:O	2.27	0.51
1:A:75:ILE:HB	1:A:92:TYR:HB3	1.94	0.50
1:B:75:ILE:HG13	1:B:152:MET:HE1	1.93	0.50
1:B:864:LEU:HD13	1:B:873:VAL:HG11	1.93	0.49
1:B:643:LYS:NZ	1:B:696:TYR:O	2.44	0.49
1:B:474:GLU:OE2	1:B:552:THR:HG22	2.11	0.49
1:B:607:TYR:CE1	1:B:614:PRO:HG2	2.47	0.49
1:A:706:PRO:HG2	1:A:709:PHE:CE2	2.47	0.49
1:A:578:SER:OG	1:A:920:GLU:HB2	2.13	0.48
1:B:799:GLU:HB3	1:B:832:VAL:HG12	1.95	0.48
1:A:128:LEU:O	2:A:1001:GOL:H32	2.14	0.47
1:A:426:ILE:HG23	1:A:456:SER:HB3	1.95	0.47
1:B:180:THR:HG23	1:B:206:SER:HB2	1.95	0.47
1:A:548:ARG:HD2	1:A:561:ASP:OD2	2.15	0.46
1:B:88:GLN:HG3	1:B:284:VAL:HG11	1.97	0.46
1:A:606:PHE:CD1	2:A:1005:GOL:H2	2.51	0.46
1:A:588:LYS:HE2	1:A:649:TYR:OH	2.16	0.46
1:B:878:TRP:CE2	1:B:906:ARG:HB2	2.52	0.45
1:B:810:LEU:HD13	1:B:814:MET:HG3	1.98	0.45
1:A:428:ILE:HG13	1:A:492:LEU:HD13	1.99	0.44
1:B:104:LYS:HE2	4:B:1509:HOH:O	2.17	0.44
1:B:805:LEU:HD22	1:B:817:MET:HE1	1.98	0.44
1:A:114:GLU:CG	1:A:659:VAL:HG11	2.48	0.44
1:B:799:GLU:HB3	1:B:832:VAL:CG1	2.47	0.44
1:B:846:MET:HE2	1:B:863:PRO:HG2	2.00	0.44
1:A:355:ALA:HB2	1:A:367:TRP:CE3	2.52	0.44
1:B:453:TYR:CZ	1:B:503:MET:HG3	2.52	0.43
1:B:525:SER:HA	1:B:916:TYR:HB2	1.99	0.43
1:B:713:ALA:N	1:B:908:MET:HG3	2.34	0.43
1:B:873:VAL:HG21	1:B:907:VAL:HG21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:491:ALA:HB2	1:A:567:LEU:HD23	2.01	0.43
1:B:353:THR:O	1:B:366:VAL:HG13	2.19	0.42
1:B:501:GLN:O	1:B:505:ARG:HG3	2.20	0.42
1:A:126:LYS:HG3	1:A:140:GLU:HG2	2.02	0.42
1:B:552:THR:HG23	4:B:1423:HOH:O	2.19	0.42
1:A:137:ILE:HD12	1:A:158:PHE:HB3	2.02	0.42
1:A:783:ARG:HD3	1:A:823:ASP:HB3	2.02	0.42
1:B:391:TRP:CH2	1:B:432:GLY:HA3	2.55	0.42
1:A:354:PRO:HD3	1:A:410:TYR:CE1	2.55	0.42
1:A:706:PRO:HG2	1:A:709:PHE:CZ	2.55	0.42
1:A:498:ASP:CG	1:A:706:PRO:HB3	2.44	0.42
1:B:631:ALA:O	1:B:635:LEU:HD23	2.21	0.41
1:B:813:SER:HB2	4:B:1166:HOH:O	2.19	0.41
1:A:224:VAL:HG12	1:A:255:GLU:HB3	2.01	0.41
1:A:789:TYR:CE2	1:A:817:MET:HG3	2.55	0.41
1:B:105:TYR:CE2	1:B:135:LEU:HD13	2.55	0.41
1:A:38:PRO:HD3	1:A:661:TYR:CZ	2.56	0.41
1:A:198:THR:OG1	1:A:295:VAL:HG11	2.21	0.41
1:A:600:ARG:HD3	1:A:626:TRP:CZ2	2.55	0.41
1:A:225:ALA:HB1	1:A:238:SER:HA	2.03	0.41
1:B:472:GLY:HA3	1:B:493:TYR:CD2	2.56	0.41
1:A:908:MET:HE3	1:A:908:MET:HB3	1.98	0.40
1:B:600:ARG:HD3	1:B:626:TRP:CZ2	2.56	0.40
1:A:529:LYS:HD2	1:A:533:TRP:CE3	2.56	0.40
1:A:829:LYS:HD2	1:A:866:GLU:OE1	2.21	0.40
1:B:788:VAL:HG22	1:B:883:ILE:HG13	2.03	0.40
1:A:249:VAL:H	2:A:1002:GOL:C3	2.17	0.40
1:B:734:PRO:HA	1:B:918:ASN:O	2.21	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	885/887 (100%)	860 (97%)	24 (3%)	1 (0%)	48	41
1	B	885/887 (100%)	862 (97%)	22 (2%)	1 (0%)	48	41
All	All	1770/1774 (100%)	1722 (97%)	46 (3%)	2 (0%)	48	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	53	ASP
1	B	53	ASP

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	754/754 (100%)	754 (100%)	0	100	100
1	B	754/754 (100%)	751 (100%)	3 (0%)	84	85
All	All	1508/1508 (100%)	1505 (100%)	3 (0%)	87	89

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

Mol	Chain	Res	Type
1	B	135	LEU
1	B	578	SER
1	B	919	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	162	ASN
1	A	173	GLN
1	A	398	GLN
1	A	543	ASN
1	A	644	GLN
1	A	666	GLN

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Mol	Chain	Res	Type
1	A	718	ASN
1	B	162	ASN
1	B	173	GLN
1	B	435	GLN
1	B	543	ASN
1	B	708	GLN
1	B	874	ASN
1	B	882	ASN
1	B	896	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 oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 17 ligands modelled in this entry, 10 are monoatomic - leaving 7 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	A	1002	-	5,5,5	0.77	0	5,5,5	1.13	1 (20%)
2	GOL	A	1001	-	5,5,5	1.04	0	5,5,5	1.26	1 (20%)
2	GOL	B	1001	-	5,5,5	1.02	0	5,5,5	0.92	0
2	GOL	A	1003	-	5,5,5	1.03	0	5,5,5	0.88	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	GOL	A	1004	-	5,5,5	1.04	0	5,5,5	1.00	0
2	GOL	B	1002	-	5,5,5	0.83	0	5,5,5	0.93	0
2	GOL	A	1005	-	5,5,5	0.88	0	5,5,5	1.05	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	A	1002	-	-	4/4/4/4	-
2	GOL	A	1001	-	-	4/4/4/4	-
2	GOL	B	1001	-	-	1/4/4/4	-
2	GOL	A	1003	-	-	0/4/4/4	-
2	GOL	A	1004	-	-	2/4/4/4	-
2	GOL	B	1002	-	-	2/4/4/4	-
2	GOL	A	1005	-	-	2/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1002	GOL	C3-C2-C1	-2.09	103.58	111.70
2	A	1001	GOL	C3-C2-C1	-2.01	103.89	111.70

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	GOL	O1-C1-C2-C3
2	A	1001	GOL	C1-C2-C3-O3
2	A	1002	GOL	O1-C1-C2-C3
2	A	1002	GOL	C1-C2-C3-O3
2	B	1002	GOL	O1-C1-C2-C3
2	A	1004	GOL	O1-C1-C2-C3
2	A	1005	GOL	C1-C2-C3-O3
2	A	1001	GOL	O1-C1-C2-O2
2	A	1001	GOL	O2-C2-C3-O3
2	A	1002	GOL	O1-C1-C2-O2
2	B	1002	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	A	1002	GOL	O2-C2-C3-O3
2	A	1004	GOL	O1-C1-C2-O2
2	A	1005	GOL	O2-C2-C3-O3
2	B	1001	GOL	O1-C1-C2-C3

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1002	GOL	2	0
2	A	1001	GOL	3	0
2	A	1005	GOL	1	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, 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	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	887/887 (100%)	0.02	8 (0%) 81 85	17, 24, 34, 53	0
1	B	887/887 (100%)	1.02	151 (17%) 4 4	15, 30, 47, 71	0
All	All	1774/1774 (100%)	0.52	159 (8%) 15 17	15, 26, 42, 71	0

All (159) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	848	ILE	5.8
1	B	863	PRO	5.5
1	B	34	THR	5.2
1	B	843	LEU	5.1
1	B	798	ALA	5.0
1	B	867	LEU	4.8
1	B	920	GLU	4.8
1	B	719	VAL	4.6
1	B	825	GLY	4.5
1	A	34	THR	4.4
1	B	713	ALA	4.3
1	B	764	LEU	4.2
1	B	757	VAL	3.8
1	B	917	GLY	3.7
1	B	720	THR	3.7
1	B	721	THR	3.7
1	B	837	TRP	3.7
1	B	861	VAL	3.6
1	B	909	THR	3.6
1	B	793	PHE	3.6
1	B	765	TYR	3.5
1	B	903	ILE	3.5
1	B	873	VAL	3.5
1	B	714	TYR	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	826	ASP	3.4
1	B	465	GLY	3.4
1	B	705	ILE	3.3
1	B	872	SER	3.3
1	B	870	THR	3.2
1	B	778	TYR	3.2
1	B	724	ALA	3.2
1	B	847	THR	3.2
1	B	916	TYR	3.1
1	B	833	SER	3.1
1	B	919	VAL	3.1
1	B	828	ALA	3.1
1	B	425	SER	3.1
1	B	449	ILE	3.1
1	B	722	THR	3.1
1	A	844	SER	3.1
1	B	801	LEU	3.1
1	B	723	VAL	3.0
1	B	716	SER	3.0
1	A	891	THR	3.0
1	B	858	ILE	3.0
1	B	896	ASN	3.0
1	B	832	VAL	3.0
1	B	853	VAL	3.0
1	B	827	PHE	3.0
1	B	179	VAL	3.0
1	B	868	SER	3.0
1	B	645	THR	3.0
1	B	533	TRP	2.9
1	B	702	ILE	2.9
1	B	831	HIS	2.8
1	B	878	TRP	2.8
1	B	509	LEU	2.8
1	B	883	ILE	2.8
1	B	452	LYS	2.8
1	B	730	TRP	2.8
1	B	635	LEU	2.8
1	B	797	THR	2.8
1	B	849	PRO	2.8
1	B	500	ALA	2.7
1	B	839	PHE	2.7
1	B	842	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	552	THR	2.7
1	B	736	TYR	2.7
1	B	821	TRP	2.7
1	B	762	ASN	2.7
1	B	771	ALA	2.6
1	A	415	TYR	2.6
1	B	844	SER	2.6
1	B	585	ASN	2.6
1	B	37	LEU	2.6
1	B	207	GLY	2.6
1	B	907	VAL	2.6
1	B	443	THR	2.6
1	B	469	TYR	2.6
1	B	709	PHE	2.6
1	B	727	LEU	2.6
1	B	835	GLY	2.6
1	B	718	ASN	2.6
1	B	570	TYR	2.5
1	B	734	PRO	2.5
1	A	917	GLY	2.5
1	A	163	ALA	2.5
1	B	753	ALA	2.5
1	B	417	SER	2.5
1	B	846	MET	2.5
1	B	851	TRP	2.5
1	B	519	ALA	2.5
1	B	755	VAL	2.5
1	B	706	PRO	2.5
1	B	816	TYR	2.4
1	B	864	LEU	2.4
1	B	813	SER	2.4
1	B	728	ALA	2.4
1	B	393	TRP	2.4
1	B	834	ASN	2.4
1	B	862	VAL	2.4
1	B	430	LEU	2.4
1	B	532	THR	2.4
1	B	733	VAL	2.4
1	B	534	SER	2.4
1	B	914	TRP	2.4
1	B	556	ALA	2.3
1	B	444	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	920	GLU	2.3
1	B	378	LEU	2.3
1	B	415	TYR	2.3
1	B	754	LYS	2.3
1	B	428	ILE	2.3
1	B	573	ILE	2.3
1	B	741	SER	2.3
1	B	879	SER	2.3
1	B	787	THR	2.3
1	B	838	THR	2.3
1	B	434	TYR	2.2
1	B	796	SER	2.2
1	A	137	ILE	2.2
1	B	766	VAL	2.2
1	B	774	HIS	2.2
1	B	874	ASN	2.2
1	B	761	ALA	2.2
1	B	763	ASN	2.2
1	B	818	VAL	2.2
1	B	464	LEU	2.2
1	B	580	ALA	2.2
1	B	470	GLY	2.2
1	B	504	ALA	2.2
1	B	461	TRP	2.1
1	B	387	ALA	2.1
1	B	631	ALA	2.1
1	B	725	ASN	2.1
1	B	790	ALA	2.1
1	B	404	PHE	2.1
1	B	830	PHE	2.1
1	B	323	THR	2.1
1	B	647	LEU	2.1
1	B	810	LEU	2.1
1	B	327	ALA	2.1
1	B	491	ALA	2.1
1	B	711	ALA	2.1
1	B	437	TYR	2.1
1	B	545	TYR	2.1
1	B	739	ARG	2.1
1	B	869	SER	2.1
1	B	576	ASN	2.1
1	B	566	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	717	ILE	2.0
1	B	649	TYR	2.0
1	B	875	THR	2.0
1	B	886	VAL	2.0
1	B	768	ILE	2.0
1	B	644	GLN	2.0
1	B	871	GLY	2.0
1	B	756	TYR	2.0
1	B	767	ARG	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, 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	B-factors(Å^2)	Q<0.9
2	GOL	B	1002	6/6	0.69	0.20	40,49,55,55	0
2	GOL	A	1004	6/6	0.75	0.24	32,42,48,56	0
2	GOL	A	1002	6/6	0.76	0.20	33,38,46,49	0
2	GOL	A	1001	6/6	0.83	0.17	19,28,37,45	0
2	GOL	A	1005	6/6	0.93	0.11	27,32,35,45	0
3	CA	A	1010	1/1	0.93	0.07	44,44,44,44	0
2	GOL	A	1003	6/6	0.95	0.10	26,28,29,31	0
2	GOL	B	1001	6/6	0.95	0.07	18,22,26,31	0
3	CA	B	1005	1/1	0.97	0.06	25,25,25,25	0
3	CA	A	1008	1/1	0.98	0.04	25,25,25,25	0
3	CA	B	1003	1/1	0.98	0.06	35,35,35,35	0
3	CA	B	1004	1/1	0.98	0.07	37,37,37,37	0
3	CA	A	1009	1/1	0.98	0.11	40,40,40,40	0
3	CA	A	1006	1/1	0.99	0.03	22,22,22,22	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	CA	A	1007	1/1	0.99	0.02	26,26,26,26	0
3	CA	B	1006	1/1	0.99	0.06	27,27,27,27	0
3	CA	B	1007	1/1	0.99	0.03	23,23,23,23	0

6.5 Other polymers (i)

There are no such residues in this entry.