



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:57 PM UTC

PDB ID : 3KLK / pdb\_00003klk  
Title : Crystal structure of Lactobacillus reuteri N-terminally truncated glucansucrase GTF180 in triclinic apo- form  
Authors : Vujicic-Zagar, A.; Pijning, T.; Kralj, S.; Eeuwema, W.; Dijkhuizen, L.; Dijkstra, B.W.  
Deposited on : 2009-11-08  
Resolution : 1.65 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](mailto: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>.

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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

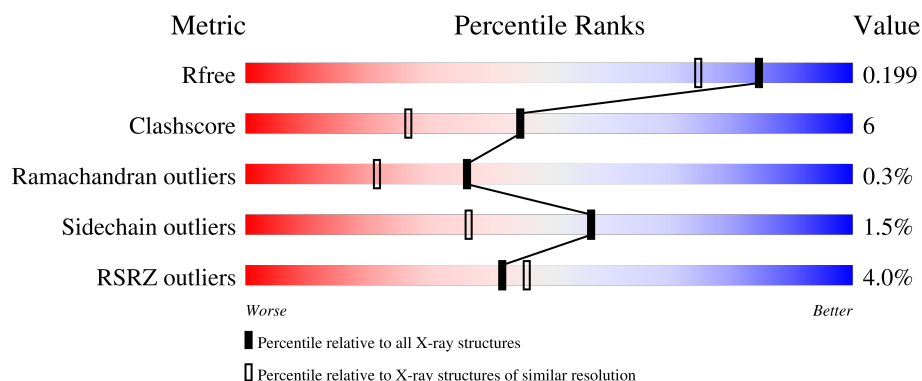
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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  $\geq 3$ , 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	1039	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	1779	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9091 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 Glucansucrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1006	Total	C	N	O	S	0	1	0
			7987	4994	1354	1618	21			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	740	MET	-	expression tag	UNP Q5SBN3
A	741	GLY	-	expression tag	UNP Q5SBN3
A	1674	LEU	PHE	SEE REMARK 999	UNP Q5SBN3
A	1773	HIS	-	expression tag	UNP Q5SBN3
A	1774	HIS	-	expression tag	UNP Q5SBN3
A	1775	HIS	-	expression tag	UNP Q5SBN3
A	1776	HIS	-	expression tag	UNP Q5SBN3
A	1777	HIS	-	expression tag	UNP Q5SBN3
A	1778	HIS	-	expression tag	UNP Q5SBN3

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

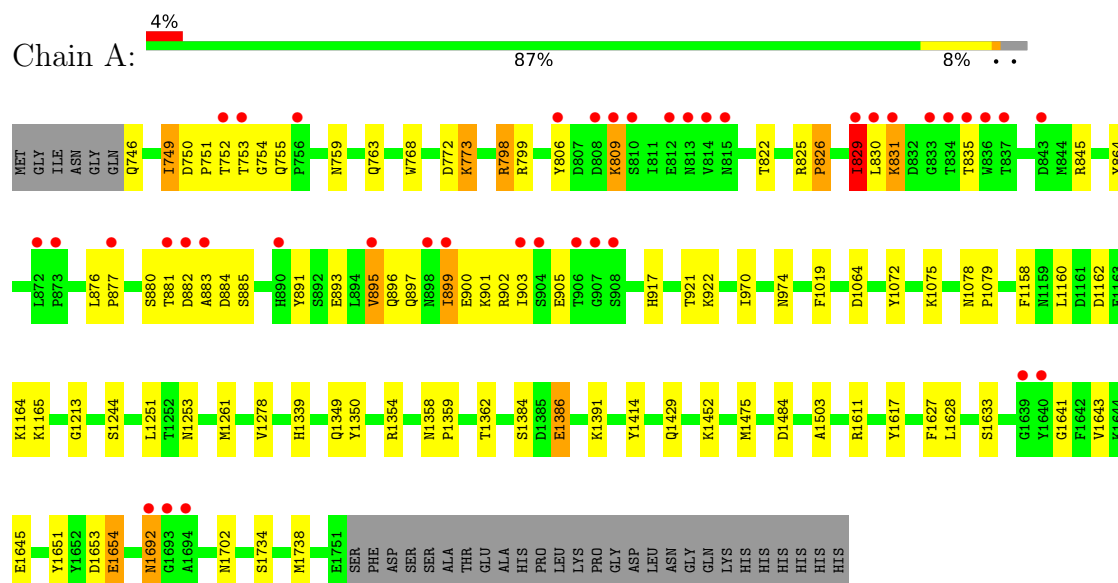
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1055	Total	O	0	0
			1055	1055		

### 3 Residue-property plots

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: Glucansucrase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.28Å 65.95Å 82.51Å 73.30° 78.46° 85.82°	Depositor
Resolution (Å)	20.00 – 1.65 20.00 – 1.65	Depositor EDS
% Data completeness (in resolution range)	96.5 (20.00-1.65) 96.4 (20.00-1.65)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 1.65Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.165 , 0.195 0.170 , 0.199	Depositor DCC
$R_{free}$ test set	6715 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.1	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 46.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	9091	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.27% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, 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	0.88	10/8170 (0.1%)	0.93	7/11114 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	893	GLU	C-N	10.66	1.47	1.33
1	A	899	ILE	C-N	7.95	1.45	1.33
1	A	809	LYS	N-CA	6.32	1.54	1.46
1	A	829	ILE	C-O	6.20	1.30	1.24
1	A	893	GLU	C-O	6.14	1.31	1.24
1	A	895	VAL	C-O	6.07	1.31	1.24
1	A	829	ILE	CA-C	6.06	1.60	1.52
1	A	809	LYS	CA-CB	6.03	1.63	1.53
1	A	826	PRO	C-O	5.77	1.31	1.23
1	A	896	GLN	C-O	5.00	1.29	1.24

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1633	SER	N-CA-C	5.92	119.69	110.17
1	A	829	ILE	N-CA-C	-5.74	100.08	108.11
1	A	893	GLU	O-C-N	5.72	127.96	122.07
1	A	826	PRO	N-CA-C	-5.63	103.34	111.22
1	A	825	ARG	N-CA-C	5.53	118.36	108.47
1	A	896	GLN	N-CA-C	-5.11	105.62	111.14
1	A	895	VAL	N-CA-C	5.04	117.34	111.05

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	7987	0	7536	88	0
2	A	1	0	0	0	0
3	A	48	0	64	11	0
4	A	1055	0	0	12	0
All	All	9091	0	7600	92	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 6.

All (92) 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)
3:A:1783:GOL:H32	4:A:171:HOH:O	1.47	1.11
1:A:826:PRO:HG2	1:A:829:ILE:CD1	1.95	0.96
1:A:806:TYR:HE2	1:A:1628:LEU:HD23	1.34	0.92
1:A:1452:LYS:NZ	3:A:1779:GOL:H31	1.87	0.90
1:A:1452:LYS:HZ3	3:A:1779:GOL:H31	1.37	0.89
1:A:1452:LYS:CE	3:A:1779:GOL:H31	2.04	0.87
1:A:826:PRO:CG	1:A:829:ILE:CD1	2.52	0.87
1:A:1452:LYS:HZ3	3:A:1779:GOL:C3	1.90	0.84
1:A:864:TYR:CE1	1:A:922:LYS:HD2	2.16	0.81
1:A:826:PRO:HG2	1:A:829:ILE:HD13	1.64	0.78
1:A:1654:GLU:H	1:A:1654:GLU:CD	1.92	0.77
1:A:902:ARG:HA	1:A:905:GLU:HG2	1.65	0.77
1:A:1452:LYS:HE2	3:A:1779:GOL:H31	1.66	0.77
1:A:773:LYS:HG2	4:A:1948:HOH:O	1.85	0.75
1:A:899:ILE:O	1:A:903:ILE:HG13	1.87	0.75
1:A:864:TYR:OH	1:A:922:LYS:HE2	1.87	0.74
1:A:864:TYR:HE1	1:A:922:LYS:HD2	1.52	0.72
1:A:749:ILE:C	1:A:749:ILE:HD12	2.14	0.72
1:A:809:LYS:HE2	4:A:2060:HOH:O	1.90	0.70
1:A:773:LYS:CG	4:A:1948:HOH:O	2.37	0.69
1:A:1452:LYS:NZ	3:A:1779:GOL:C3	2.52	0.69
1:A:806:TYR:HE2	1:A:1628:LEU:CD2	2.05	0.68
1:A:884:ASP:OD1	1:A:885:SER:N	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:806:TYR:CE2	1:A:1628:LEU:HD23	2.24	0.67
1:A:826:PRO:HG3	1:A:829:ILE:CD1	2.25	0.65
1:A:845:ARG:NE	1:A:900:GLU:OE2	2.22	0.64
1:A:826:PRO:CG	1:A:829:ILE:HD11	2.27	0.64
1:A:798:ARG:HG3	1:A:799:ARG:N	2.13	0.64
1:A:1075:LYS:HD3	4:A:1858:HOH:O	1.99	0.61
1:A:753:THR:C	1:A:755:GLN:H	2.08	0.61
1:A:806:TYR:CE2	1:A:1628:LEU:CD2	2.85	0.60
1:A:1627:PHE:C	1:A:1628:LEU:HD22	2.27	0.59
1:A:895:VAL:O	1:A:899:ILE:HG13	2.02	0.59
1:A:864:TYR:HE1	1:A:922:LYS:CD	2.15	0.58
1:A:749:ILE:HD12	1:A:750:ASP:C	2.28	0.57
1:A:1339:HIS:HD2	4:A:235:HOH:O	1.87	0.57
1:A:1261:MET:HE2	1:A:1261:MET:HA	1.87	0.56
1:A:826:PRO:HG3	1:A:829:ILE:HD11	1.87	0.56
1:A:1213:GLY:HA2	1:A:1414:TYR:CE2	2.40	0.55
1:A:897:GLN:O	1:A:901:LYS:HG3	2.07	0.55
1:A:753:THR:O	1:A:755:GLN:HG2	2.07	0.54
1:A:759:ASN:HB2	1:A:773:LYS:HG3	1.89	0.54
1:A:1475:MET:HG3	1:A:1484:ASP:HB3	1.91	0.53
1:A:753:THR:C	1:A:755:GLN:N	2.65	0.52
3:A:1783:GOL:C3	4:A:171:HOH:O	2.27	0.52
1:A:753:THR:O	1:A:755:GLN:N	2.42	0.52
1:A:1452:LYS:HZ3	3:A:1779:GOL:C2	2.22	0.52
1:A:1611:ARG:HG2	1:A:1617:TYR:CD1	2.46	0.51
1:A:1160:LEU:HG	1:A:1164:LYS:HE3	1.92	0.51
1:A:830:LEU:HB3	1:A:900:GLU:HG3	1.94	0.50
1:A:884:ASP:OD1	1:A:884:ASP:C	2.55	0.49
1:A:831:LYS:CB	1:A:835:THR:O	2.61	0.49
1:A:1253:ASN:HB3	1:A:1278:VAL:HB	1.95	0.49
1:A:826:PRO:CG	1:A:829:ILE:HD12	2.37	0.49
1:A:876:LEU:HD21	1:A:891:TYR:CE1	2.48	0.49
1:A:880:SER:O	1:A:883:ALA:N	2.45	0.49
1:A:880:SER:C	1:A:882:ASP:N	2.70	0.48
1:A:831:LYS:HB3	1:A:835:THR:O	2.14	0.48
1:A:1654:GLU:CD	1:A:1654:GLU:N	2.62	0.47
1:A:902:ARG:CA	1:A:905:GLU:HG2	2.39	0.47
1:A:763:GLN:HB2	1:A:768:TRP:CH2	2.51	0.46
1:A:880:SER:O	1:A:882:ASP:N	2.49	0.46
1:A:1653:ASP:OD1	1:A:1653:ASP:C	2.57	0.46
1:A:1019:PHE:CG	1:A:1503:ALA:HB2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1429:GLN:NE2	4:A:677:HOH:O	2.49	0.45
1:A:1165:LYS:HD2	4:A:701:HOH:O	2.16	0.45
1:A:1358:ASN:HB2	1:A:1359:PRO:CD	2.47	0.45
1:A:1349:GLN:HG2	1:A:1350:TYR:CD2	2.52	0.45
1:A:1643:VAL:HG12	1:A:1645:GLU:HG3	1.99	0.44
1:A:752:THR:C	1:A:753:THR:HG23	2.43	0.44
3:A:1779:GOL:O3	3:A:1779:GOL:O1	2.26	0.44
3:A:1783:GOL:C1	4:A:171:HOH:O	2.65	0.44
1:A:880:SER:C	1:A:882:ASP:H	2.26	0.44
1:A:970:ILE:HD11	1:A:1072:TYR:CE1	2.53	0.43
1:A:1158:PHE:HB2	1:A:1162:ASP:HB2	2.01	0.43
1:A:763:GLN:HB2	1:A:768:TRP:CZ3	2.53	0.42
1:A:752:THR:C	1:A:753:THR:CG2	2.93	0.42
1:A:974:ASN:HB3	4:A:2093:HOH:O	2.19	0.42
1:A:1078:ASN:N	1:A:1079:PRO:CD	2.82	0.42
1:A:749:ILE:HD12	1:A:751:PRO:N	2.35	0.42
1:A:1641:GLY:HA2	1:A:1651:TYR:CE1	2.54	0.42
1:A:746:GLN:NE2	4:A:2097:HOH:O	2.53	0.41
1:A:772:ASP:C	1:A:772:ASP:OD1	2.63	0.41
1:A:917:HIS:O	1:A:921:THR:HG23	2.21	0.41
1:A:876:LEU:HA	1:A:877:PRO:HD3	1.91	0.41
1:A:830:LEU:O	1:A:831:LYS:C	2.63	0.41
1:A:1734:SER:HA	1:A:1738:MET:O	2.21	0.41
1:A:902:ARG:O	1:A:905:GLU:HG2	2.20	0.41
1:A:1384:SER:OG	1:A:1386:GLU:HB2	2.21	0.41
1:A:1244:SER:O	1:A:1251:LEU:HD12	2.21	0.40
1:A:902:ARG:HA	1:A:905:GLU:CG	2.42	0.40
1:A:1064:ASP:CG	1:A:1064:ASP:O	2.64	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	1006/1039 (97%)	980 (97%)	23 (2%)	3 (0%)	36	21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	881	THR
1	A	1692	ASN
1	A	754	GLY

### 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	858/883 (97%)	845 (98%)	13 (2%)	57	37

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	749	ILE
1	A	773	LYS
1	A	798	ARG
1	A	822	THR
1	A	829	ILE
1	A	831	LYS
1	A	1354	ARG
1	A	1362	THR
1	A	1386	GLU
1	A	1391	LYS
1	A	1654	GLU
1	A	1692	ASN
1	A	1702	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	746	GLN
1	A	755	GLN
1	A	867	GLN
1	A	1074	ASN
1	A	1078	ASN
1	A	1145	GLN
1	A	1148	GLN
1	A	1174	GLN
1	A	1182	ASN
1	A	1573	GLN
1	A	1648	ASN
1	A	1650	HIS
1	A	1679	ASN
1	A	1731	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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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	$\# Z  > 2$
3	GOL	A	1779	-	5,5,5	0.25	0	5,5,5	0.90	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	GOL	A	1786	-	5,5,5	0.31	0	5,5,5	0.66	0
3	GOL	A	1784	-	5,5,5	0.41	0	5,5,5	0.87	0
3	GOL	A	1781	-	5,5,5	0.32	0	5,5,5	0.67	0
3	GOL	A	1780	-	5,5,5	0.28	0	5,5,5	0.64	0
3	GOL	A	1785	-	5,5,5	0.17	0	5,5,5	0.81	0
3	GOL	A	1782	-	5,5,5	0.39	0	5,5,5	0.25	0
3	GOL	A	1783	-	5,5,5	0.44	0	5,5,5	0.44	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
3	GOL	A	1779	-	-	0/4/4/4	-
3	GOL	A	1786	-	-	2/4/4/4	-
3	GOL	A	1784	-	-	0/4/4/4	-
3	GOL	A	1781	-	-	1/4/4/4	-
3	GOL	A	1780	-	-	2/4/4/4	-
3	GOL	A	1785	-	-	0/4/4/4	-
3	GOL	A	1782	-	-	0/4/4/4	-
3	GOL	A	1783	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1783	GOL	C1-C2-C3-O3
3	A	1783	GOL	O2-C2-C3-O3
3	A	1780	GOL	O1-C1-C2-C3
3	A	1786	GOL	C1-C2-C3-O3
3	A	1786	GOL	O2-C2-C3-O3
3	A	1780	GOL	O1-C1-C2-O2
3	A	1783	GOL	O1-C1-C2-O2
3	A	1781	GOL	C1-C2-C3-O3
3	A	1783	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1779	GOL	8	0
3	A	1783	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	1006/1039 (96%)	-0.44	40 (3%)	42 46	6, 18, 50, 79	1 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	814	VAL	5.9
1	A	815	ASN	5.7
1	A	837	THR	4.8
1	A	813	ASN	4.5
1	A	836	TRP	4.2
1	A	1692	ASN	4.0
1	A	903	ILE	3.8
1	A	831	LYS	3.6
1	A	808	ASP	3.5
1	A	899	ILE	3.4
1	A	835	THR	3.4
1	A	830	LEU	3.4
1	A	881	THR	3.4
1	A	1693	GLY	3.2
1	A	1639	GLY	3.0
1	A	882	ASP	3.0
1	A	907	GLY	2.9
1	A	883	ALA	2.7
1	A	753	THR	2.7
1	A	904	SER	2.6
1	A	829	ILE	2.6
1	A	1694	ALA	2.6
1	A	833	GLY	2.5
1	A	895	VAL	2.5
1	A	877	PRO	2.5
1	A	810	SER	2.5
1	A	752	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	809	LYS	2.3
1	A	806	TYR	2.3
1	A	756	PRO	2.3
1	A	873	PRO	2.3
1	A	906	THR	2.3
1	A	812	GLU	2.2
1	A	843	ASP	2.2
1	A	890	HIS	2.2
1	A	898	ASN	2.1
1	A	834	THR	2.1
1	A	908	SER	2.0
1	A	872	LEU	2.0
1	A	1640	TYR	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<sup>th</sup> 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
3	GOL	A	1786	6/6	0.76	0.14	29,37,42,47	0
3	GOL	A	1779	6/6	0.78	0.13	25,31,41,42	0
3	GOL	A	1783	6/6	0.86	0.10	27,29,33,45	0
3	GOL	A	1780	6/6	0.90	0.12	18,20,41,42	0
3	GOL	A	1785	6/6	0.91	0.08	28,29,31,34	0
3	GOL	A	1782	6/6	0.95	0.06	18,22,24,26	0
3	GOL	A	1784	6/6	0.96	0.06	17,20,23,34	0
3	GOL	A	1781	6/6	0.97	0.04	10,10,11,12	0
2	CA	A	1	1/1	0.99	0.02	9,9,9,9	0



## 6.5 Other polymers [i](#)

There are no such residues in this entry.