



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 2WAN / pdb_00002wan
Title : Pullulanase from *Bacillus acidopullulyticus*
Authors : Turkenburg, J.P.; Brzozowski, A.M.; Svendsen, A.; Borchert, T.V.; Davies, G.J.; Wilson, K.S.
Deposited on : 2009-02-10
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

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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.42

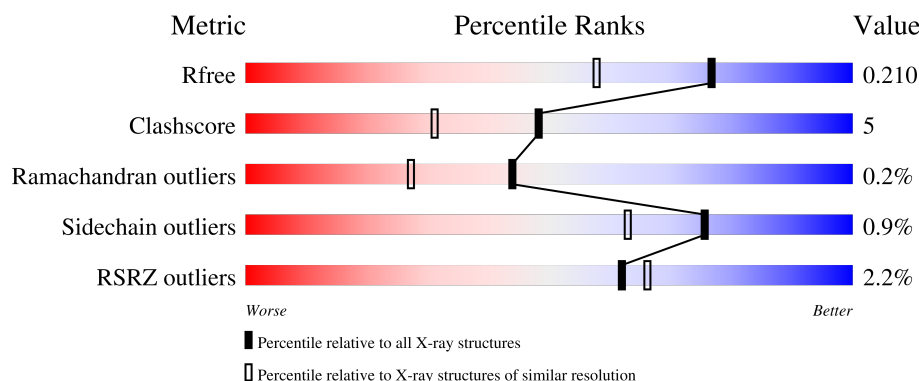
1 Overall quality at a glance

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}	164625	2328 (1.66-1.66)
Clashscore	180529	2515 (1.66-1.66)
Ramachandran outliers	177936	2475 (1.66-1.66)
Sidechain outliers	177891	2475 (1.66-1.66)
RSRZ outliers	164620	2328 (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 ≥ 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	921	

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
2	GOL	A	1929	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACT	A	1934	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7477 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 PULLULANASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	810	Total	C	N	O	S	0	23	0
			6416	4041	1077	1277	21			

- 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		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
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	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	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		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	977	Total	O	0	0
			977	977		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	163.43Å 61.98Å 109.67Å 90.00° 109.10° 90.00°	Depositor
Resolution (Å)	105.41 – 1.65 103.63 – 1.65	Depositor EDS
% Data completeness (in resolution range)	95.1 (105.41-1.65) 95.1 (103.63-1.65)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.5.0070	Depositor
R, R_{free}	0.175 , 0.210 0.174 , 0.210	Depositor DCC
R_{free} test set	6035 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	16.2	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 50.3	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	7477	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.54	1/6617 (0.0%)	0.60	0/9014

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	145	ILE	CA-CB	-5.87	1.41	1.54

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	6416	0	6227	66	0
2	A	48	0	61	12	0
3	A	20	0	0	0	0
4	A	16	0	12	2	0
5	A	977	0	0	10	0
All	All	7477	0	6300	69	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 5.

All (69) 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:386[A]:GLY:O	1:A:387[A]:THR:CG2	1.72	1.37
1:A:386[A]:GLY:O	1:A:387[A]:THR:HG22	1.33	1.13
1:A:479[D]:LYS:HA	1:A:479[D]:LYS:HE2	1.34	1.10
1:A:384[B]:VAL:HG12	1:A:385[B]:ASN:H	1.00	1.10
1:A:386[A]:GLY:O	1:A:387[A]:THR:HG23	1.52	1.09
1:A:384[B]:VAL:O	1:A:386[B]:GLY:N	1.92	1.03
1:A:209:THR:HG22	1:A:232:THR:HG22	1.42	0.99
1:A:384[B]:VAL:CG1	1:A:385[B]:ASN:H	1.76	0.96
1:A:384[B]:VAL:HG12	1:A:385[B]:ASN:N	1.80	0.94
1:A:386[A]:GLY:C	1:A:387[A]:THR:HG23	1.91	0.89
1:A:543:GLN:HE21	2:A:1927:GOL:H31	1.37	0.88
1:A:173:GLN:HE22	1:A:183:SER:H	1.23	0.85
1:A:402:ASN:HD21	1:A:499:GLU:H	1.23	0.82
1:A:878:SER:H	2:A:1925:GOL:H31	1.43	0.81
1:A:612:ASN:HB3	5:A:2598:HOH:O	1.89	0.73
1:A:572:TYR:O	1:A:596:HIS:HD2	1.73	0.72
1:A:711:LYS:HE2	4:A:1934:ACT:H3	1.71	0.72
1:A:512:ARG:HE	1:A:513:ASN:HD21	1.37	0.71
1:A:437:GLU:OE2	1:A:735:HIS:HD2	1.74	0.70
1:A:479[C]:LYS:HD3	1:A:544:GLN:HE21	1.56	0.70
1:A:557:THR:H	1:A:590:ASN:HD21	1.39	0.69
1:A:542:HIS:HE1	1:A:617:ASP:OD2	1.76	0.69
1:A:690:ASN:HD22	1:A:692:PHE:H	1.42	0.67
1:A:382:VAL:HG12	1:A:384[B]:VAL:HG23	1.78	0.65
1:A:690:ASN:ND2	1:A:692:PHE:H	1.95	0.65
1:A:900:LYS:HA	2:A:1929:GOL:H12	1.80	0.62
1:A:146:ASP:OD2	1:A:286:HIS:HD2	1.82	0.61
1:A:878:SER:N	2:A:1925:GOL:H31	2.15	0.61
1:A:818:SER:O	1:A:822:HIS:HD2	1.85	0.60
1:A:402:ASN:ND2	1:A:499:GLU:H	1.96	0.60
1:A:479[D]:LYS:HA	1:A:479[D]:LYS:CE	2.20	0.60
1:A:493:GLU:HG2	1:A:513:ASN:HD22	1.67	0.58
1:A:640:HIS:HE1	1:A:646:ILE:O	1.85	0.58
1:A:688:ASP:OD2	2:A:1922:GOL:H12	2.04	0.57
1:A:901:SER:HB2	2:A:1928:GOL:H2	1.88	0.56
1:A:512:ARG:HE	1:A:513:ASN:ND2	2.05	0.54
1:A:642:ILE:HD13	5:A:2589:HOH:O	2.07	0.53
2:A:1928:GOL:O2	2:A:1929:GOL:H32	2.09	0.52
1:A:479[C]:LYS:HD3	1:A:544:GLN:NE2	2.24	0.52
1:A:555:ASN:OD1	1:A:556:HIS:HD2	1.93	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:822:HIS:HE1	5:A:2462:HOH:O	1.93	0.51
1:A:304:ASN:OD1	1:A:596:HIS:HE1	1.94	0.50
1:A:735:HIS:HE1	5:A:2731:HOH:O	1.96	0.49
1:A:761:GLU:OE1	2:A:1924:GOL:H31	2.14	0.48
1:A:479[D]:LYS:HD2	1:A:544:GLN:HB3	1.96	0.48
2:A:1922:GOL:H11	5:A:2953:HOH:O	2.12	0.47
1:A:507:TRP:CD2	1:A:793:TYR:HA	2.49	0.47
1:A:901:SER:H	2:A:1929:GOL:H12	1.80	0.47
1:A:479[D]:LYS:HE2	1:A:479[D]:LYS:CA	2.24	0.47
1:A:185:ASP:HA	1:A:218:HIS:CD2	2.49	0.47
1:A:556:HIS:HE1	5:A:2531:HOH:O	1.99	0.46
1:A:174:GLN:H	1:A:181:ASN:HD21	1.64	0.46
1:A:286:HIS:HE1	5:A:2078:HOH:O	1.98	0.45
1:A:186:ASP:OD1	1:A:188:HIS:HB2	2.16	0.45
1:A:174:GLN:H	1:A:181:ASN:ND2	2.16	0.44
1:A:184:PRO:HG3	1:A:220:TRP:CE2	2.52	0.44
1:A:901:SER:H	2:A:1929:GOL:C1	2.31	0.44
1:A:437:GLU:OE2	1:A:735:HIS:CD2	2.63	0.44
1:A:732:VAL:O	2:A:1922:GOL:C1	2.67	0.43
1:A:764:HIS:HE1	5:A:2726:HOH:O	2.00	0.43
1:A:911:VAL:HG22	1:A:917:LEU:HD22	2.01	0.42
1:A:480:GLU:HG2	1:A:814:PHE:HE1	1.85	0.42
1:A:173:GLN:NE2	1:A:183:SER:H	2.03	0.42
1:A:818:SER:O	1:A:822:HIS:CD2	2.71	0.42
1:A:147:LYS:HE3	1:A:287:ASN:ND2	2.36	0.41
1:A:384[B]:VAL:CG1	1:A:385[B]:ASN:N	2.49	0.41
1:A:306:LEU:HB2	1:A:572:TYR:CE1	2.56	0.41
4:A:1934:ACT:H1	5:A:2698:HOH:O	2.21	0.41
1:A:896:GLN:HG2	5:A:2923:HOH:O	2.21	0.40

There are no symmetry-related clashes.

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	832/921 (90%)	799 (96%)	29 (4%)	4 (0%)	25 11

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	385[A]	ASN
1	A	385[B]	ASN
1	A	384[A]	VAL
1	A	384[B]	VAL

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	715/785 (91%)	709 (99%)	6 (1%)	79 68

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	164	VAL
1	A	172	LEU
1	A	205	LEU
1	A	221	ASN
1	A	620	ARG
1	A	920	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	GLN
1	A	174	GLN
1	A	181	ASN
1	A	200	GLN
1	A	211	GLN
1	A	228	ASN

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Mol	Chain	Res	Type
1	A	286	HIS
1	A	354	GLN
1	A	358	GLN
1	A	389	GLN
1	A	402	ASN
1	A	513	ASN
1	A	542	HIS
1	A	543	GLN
1	A	544	GLN
1	A	556	HIS
1	A	590	ASN
1	A	596	HIS
1	A	640	HIS
1	A	690	ASN
1	A	719	GLN
1	A	735	HIS
1	A	764	HIS
1	A	822	HIS
1	A	826	GLN
1	A	908	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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	$\# Z > 2$
1	CSX	A	588	1	3,6,7	0.65	0	1,6,8	0.04	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
1	CSX	A	588	1	-	0/1/5/7	-

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

16 ligands are modelled in this entry.

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	$\# Z > 2$
4	ACT	A	1934	-	3,3,3	0.80	0	3,3,3	0.93	0
4	ACT	A	1935	-	3,3,3	0.81	0	3,3,3	0.86	0
4	ACT	A	1936	-	3,3,3	0.85	0	3,3,3	1.10	0
2	GOL	A	1923	-	5,5,5	0.25	0	5,5,5	0.29	0
2	GOL	A	1927	-	5,5,5	0.32	0	5,5,5	0.36	0
3	SO4	A	1932	-	4,4,4	0.21	0	6,6,6	0.18	0
3	SO4	A	1933	-	4,4,4	0.14	0	6,6,6	0.09	0
4	ACT	A	1937	-	3,3,3	0.77	0	3,3,3	1.37	0
2	GOL	A	1926	-	5,5,5	0.45	0	5,5,5	0.67	0
2	GOL	A	1925	-	5,5,5	0.38	0	5,5,5	0.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	A	1929	2	5,5,5	0.40	0	5,5,5	0.45	0
2	GOL	A	1922	-	5,5,5	0.36	0	5,5,5	0.80	0
2	GOL	A	1924	-	5,5,5	0.41	0	5,5,5	0.70	0
2	GOL	A	1928	2	5,5,5	0.41	0	5,5,5	0.24	0
3	SO4	A	1931	-	4,4,4	0.17	0	6,6,6	0.32	0
3	SO4	A	1930	-	4,4,4	0.17	0	6,6,6	0.39	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	1923	-	-	0/4/4/4	-
2	GOL	A	1927	-	-	0/4/4/4	-
2	GOL	A	1926	-	-	4/4/4/4	-
2	GOL	A	1925	-	-	0/4/4/4	-
2	GOL	A	1929	2	-	2/4/4/4	-
2	GOL	A	1922	-	-	0/4/4/4	-
2	GOL	A	1924	-	-	3/4/4/4	-
2	GOL	A	1928	2	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1924	GOL	C1-C2-C3-O3
2	A	1926	GOL	C1-C2-C3-O3
2	A	1928	GOL	O1-C1-C2-O2
2	A	1928	GOL	O1-C1-C2-C3
2	A	1928	GOL	C1-C2-C3-O3
2	A	1929	GOL	O1-C1-C2-O2
2	A	1929	GOL	O1-C1-C2-C3
2	A	1926	GOL	O1-C1-C2-C3
2	A	1924	GOL	O2-C2-C3-O3
2	A	1926	GOL	O2-C2-C3-O3
2	A	1928	GOL	O2-C2-C3-O3
2	A	1924	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	A	1926	GOL	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1934	ACT	2	0
2	A	1927	GOL	1	0
2	A	1925	GOL	2	0
2	A	1929	GOL	4	0
2	A	1922	GOL	3	0
2	A	1924	GOL	1	0
2	A	1928	GOL	2	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, 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	809/921 (87%)	-0.37	18 (2%) 62 66	4, 15, 41, 62	23 (2%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	385[A]	ASN	3.8
1	A	386[A]	GLY	3.8
1	A	261	ALA	3.7
1	A	112	PRO	3.6
1	A	259	ASN	3.6
1	A	264	THR	3.0
1	A	145	ILE	3.0
1	A	692	PHE	2.7
1	A	263	PRO	2.6
1	A	260	GLN	2.5
1	A	137	ASP	2.4
1	A	188	HIS	2.4
1	A	262	PHE	2.3
1	A	153	ILE	2.2
1	A	258	PRO	2.2
1	A	237	GLY	2.2
1	A	236	GLY	2.1
1	A	233	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(\AA^2)	Q<0.9
1	CSX	A	588	7/8	0.97	0.05	18,21,32,33	0

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, 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(\AA^2)	Q<0.9
2	GOL	A	1925	6/6	0.71	0.24	42,44,51,52	0
4	ACT	A	1936	4/4	0.77	0.16	36,39,41,42	0
2	GOL	A	1927	6/6	0.80	0.16	34,44,48,51	0
2	GOL	A	1928	6/6	0.81	0.21	18,21,25,32	6
2	GOL	A	1926	6/6	0.82	0.19	28,39,42,48	0
4	ACT	A	1934	4/4	0.83	0.17	27,29,31,32	0
2	GOL	A	1929	6/6	0.84	0.20	15,23,27,33	6
4	ACT	A	1937	4/4	0.87	0.15	44,44,45,45	0
2	GOL	A	1922	6/6	0.88	0.18	10,20,34,41	0
3	SO4	A	1933	5/5	0.90	0.10	56,56,58,59	0
2	GOL	A	1924	6/6	0.92	0.13	15,27,31,34	0
3	SO4	A	1930	5/5	0.94	0.14	20,25,29,29	0
3	SO4	A	1932	5/5	0.95	0.15	28,36,40,46	0
4	ACT	A	1935	4/4	0.95	0.11	25,29,29,30	0
2	GOL	A	1923	6/6	0.96	0.06	17,19,20,20	0
3	SO4	A	1931	5/5	0.96	0.13	22,31,33,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.