



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:22 PM UTC

PDB ID : 2W18 / pdb_00002w18
Title : Crystal structure of the C-terminal WD40 domain of human PALB2
Authors : Oliver, A.W.; Pearl, L.H.
Deposited on : 2008-10-15
Resolution : 1.90 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

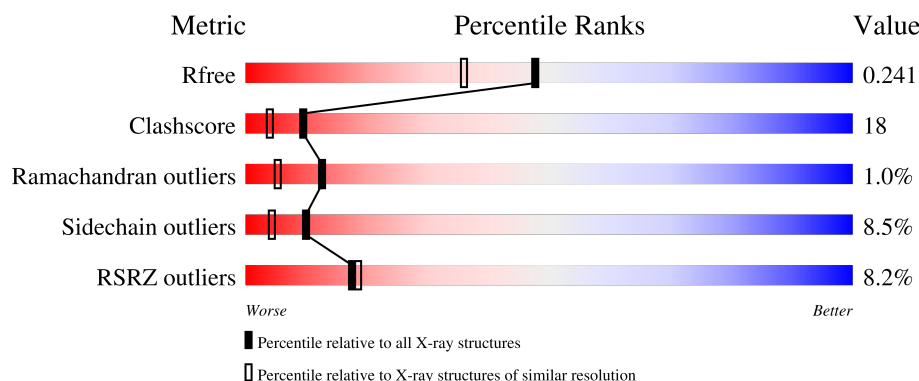
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	356	7% 58% 24% • 14%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2543 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 PARTNER AND LOCALIZER OF BRCA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	306	2326	1510	371	427	18	0	2	0

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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		

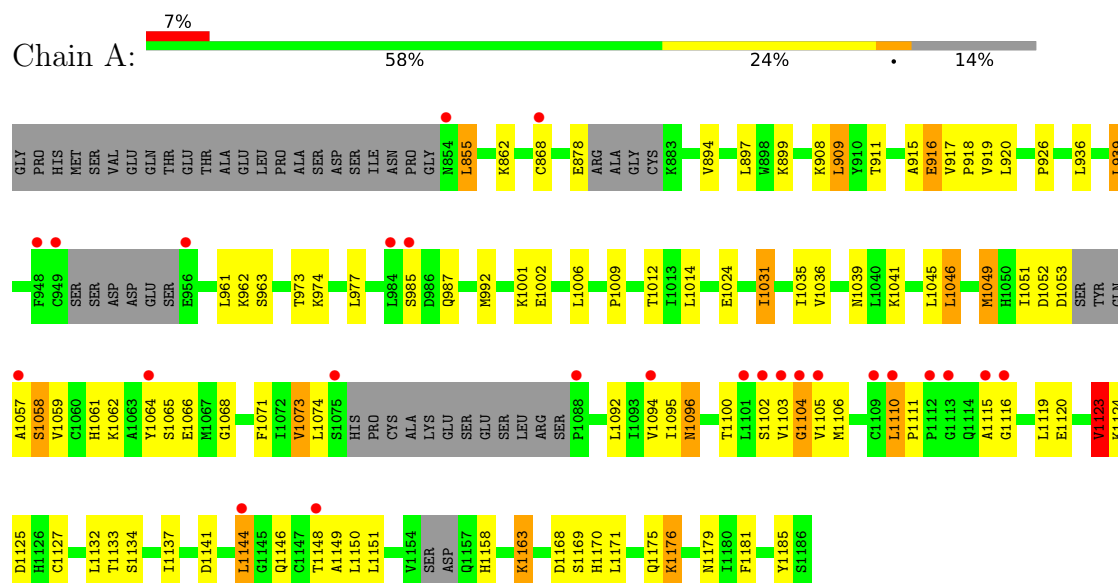
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	187	Total 187	O 187	0	0

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: PARTNER AND LOCALIZER OF BRCA2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	80.78Å 64.65Å 77.25Å 90.00° 109.65° 90.00°	Depositor
Resolution (Å)	49.27 – 1.90 49.27 – 1.90	Depositor EDS
% Data completeness (in resolution range)	93.5 (49.27-1.90) 93.8 (49.27-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 1.90Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.209 , 0.246 0.204 , 0.241	Depositor DCC
R_{free} test set	1405 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 55.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2543	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.33% 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

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	0.70	0/2381	0.94	7/3243 (0.2%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1110	LEU	CA-C-N	6.53	127.10	120.38
1	A	1110	LEU	C-N-CA	6.53	127.10	120.38
1	A	1125	ASP	CB-CA-C	-6.13	108.89	117.23
1	A	1104	GLY	N-CA-C	-5.56	104.85	111.36
1	A	1096	ASN	CA-C-N	5.33	125.00	119.56
1	A	1096	ASN	C-N-CA	5.33	125.00	119.56
1	A	1123	VAL	N-CA-C	5.29	116.34	108.46

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	2326	0	2293	81	0
2	A	30	0	40	6	0
3	A	187	0	0	10	0
All	All	2543	0	2333	84	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 18.

All (84) 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:1057:ALA:HA	1:A:1058:SER:HB2	1.30	1.04
1:A:961:LEU:HD13	1:A:992:MET:HE1	1.36	1.04
1:A:1057:ALA:HA	1:A:1058:SER:CB	1.93	0.98
1:A:1014:LEU:HD21	1:A:1031:ILE:HG22	1.48	0.95
2:A:2188:GOL:H11	3:A:3182:HOH:O	1.66	0.92
1:A:1035:ILE:HG23	1:A:1051:ILE:HD11	1.53	0.91
1:A:1049:MET:HG3	1:A:1100:THR:HG23	1.55	0.88
2:A:2190:GOL:H12	3:A:3082:HOH:O	1.74	0.87
1:A:1057:ALA:CA	1:A:1058:SER:HB2	2.10	0.81
1:A:1062:LYS:HB3	1:A:1073:VAL:HG13	1.67	0.76
1:A:855:LEU:HD13	1:A:1185:TYR:HD2	1.56	0.70
1:A:1103:VAL:HB	3:A:3144:HOH:O	1.91	0.70
1:A:1141:ASP:HB3	1:A:1144:LEU:HB2	1.74	0.69
1:A:1175:GLN:HE21	1:A:1181:PHE:HE2	1.40	0.69
2:A:2188:GOL:H12	3:A:3135:HOH:O	1.95	0.66
1:A:1124:LYS:HD3	2:A:2191:GOL:H11	1.78	0.66
1:A:1119:LEU:HD22	1:A:1133:THR:HA	1.79	0.65
1:A:1001:LYS:HG2	1:A:1002:GLU:HG2	1.77	0.65
1:A:1062:LYS:HE3	1:A:1064:TYR:CE1	2.33	0.62
1:A:868:CYS:SG	1:A:1176:LYS:HG2	2.40	0.62
1:A:1127:CYS:HB2	3:A:3187:HOH:O	1.99	0.62
1:A:862:LYS:HA	1:A:1179:ASN:OD1	2.00	0.61
1:A:1168:ASP:CG	1:A:1170:HIS:HD1	2.09	0.60
1:A:1035:ILE:HD11	1:A:1049:MET:HE3	1.84	0.59
1:A:1035:ILE:CG2	1:A:1051:ILE:HD11	2.31	0.59
1:A:919:VAL:HG22	1:A:936:LEU:CD2	2.33	0.59
1:A:1062:LYS:HG2	1:A:1064:TYR:CZ	2.37	0.59
1:A:1071:PHE:CD1	1:A:1094:VAL:HG22	2.39	0.58
1:A:1092:LEU:HB3	1:A:1106:MET:HB2	1.85	0.58
1:A:894:VAL:CG2	1:A:911:THR:HG23	2.34	0.57
1:A:1012:THR:HB	1:A:1031:ILE:HD12	1.88	0.55
1:A:1039:ASN:HB2	1:A:1046:LEU:HD11	1.87	0.54
1:A:1061:HIS:HB3	1:A:1123:VAL:HG13	1.89	0.54
1:A:899:LYS:HB2	1:A:909:LEU:HD22	1.90	0.54
1:A:908:LYS:HD3	1:A:911:THR:OG1	2.08	0.54
1:A:1095:ILE:HG12	1:A:1102:SER:CB	2.39	0.53
1:A:894:VAL:HG23	1:A:911:THR:HG23	1.91	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1100:THR:O	1:A:1100:THR:HG22	2.10	0.52
1:A:1095:ILE:HG12	1:A:1102:SER:HB2	1.92	0.51
1:A:1111:PRO:HG2	1:A:1150:LEU:HD21	1.92	0.51
1:A:961:LEU:HD13	1:A:992:MET:CE	2.25	0.50
1:A:1062:LYS:HG2	1:A:1064:TYR:OH	2.11	0.50
1:A:1110:LEU:HD21	1:A:1132:LEU:HD11	1.92	0.50
1:A:1168:ASP:OD1	1:A:1170:HIS:ND1	2.31	0.50
1:A:1051:ILE:O	1:A:1052:ASP:C	2.54	0.49
1:A:915:ALA:O	1:A:916:GLU:CB	2.61	0.48
1:A:1103:VAL:HG12	1:A:1104:GLY:O	2.13	0.48
1:A:1169:SER:HA	2:A:2191:GOL:H12	1.96	0.48
1:A:1134:SER:O	1:A:1158:HIS:CE1	2.66	0.48
1:A:1137:ILE:HB	1:A:1151:LEU:HB3	1.95	0.48
1:A:868:CYS:SG	1:A:1176:LYS:HD3	2.54	0.48
1:A:1144:LEU:HB3	1:A:1146:GLN:HG2	1.97	0.47
1:A:963:SER:HB3	2:A:2190:GOL:O3	2.13	0.47
1:A:973:THR:O	1:A:974:LYS:HB2	2.15	0.47
1:A:1148:THR:O	1:A:1149:ALA:HB2	2.16	0.46
1:A:919:VAL:HG22	1:A:936:LEU:HD22	1.99	0.45
1:A:1024:GLU:HB2	1:A:1041:LYS:HE3	1.98	0.45
1:A:1119:LEU:HD13	1:A:1133:THR:HG22	1.99	0.44
1:A:926:PRO:HB2	1:A:973:THR:HA	2.00	0.44
1:A:987:GLN:OE1	1:A:1009:PRO:HD2	2.17	0.43
1:A:1049:MET:HE2	1:A:1049:MET:HB2	1.73	0.43
1:A:897:LEU:HD23	1:A:909:LEU:HB2	2.01	0.43
1:A:1041:LYS:HE3	3:A:3126:HOH:O	2.19	0.43
1:A:1062:LYS:CG	1:A:1064:TYR:CZ	3.01	0.43
1:A:961:LEU:CD1	1:A:992:MET:HE1	2.26	0.43
1:A:1039:ASN:HB2	1:A:1046:LEU:CD1	2.48	0.43
1:A:1110:LEU:HD11	1:A:1132:LEU:HD13	2.01	0.43
1:A:1163:LYS:NZ	3:A:3150:HOH:O	2.51	0.43
1:A:1036:VAL:HB	1:A:1045:LEU:HD11	2.00	0.43
1:A:1065:SER:O	1:A:1066:GLU:HG3	2.19	0.42
1:A:1103:VAL:O	1:A:1105:VAL:HG23	2.20	0.42
1:A:1175:GLN:NE2	1:A:1181:PHE:HE2	2.14	0.42
1:A:1105:VAL:HG12	1:A:1105:VAL:O	2.19	0.42
1:A:1096:ASN:C	1:A:1096:ASN:HD22	2.28	0.41
1:A:1127:CYS:CB	3:A:3187:HOH:O	2.65	0.41
1:A:1115:ALA:HA	1:A:1116:GLY:HA2	1.83	0.41
1:A:1096:ASN:ND2	3:A:3142:HOH:O	2.52	0.41
1:A:1119:LEU:HD12	1:A:1119:LEU:HA	1.78	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:917:VAL:HA	1:A:918:PRO:HD3	1.80	0.40
1:A:1096:ASN:O	1:A:1100:THR:N	2.53	0.40
1:A:1141:ASP:CG	1:A:1144:LEU:HD22	2.46	0.40
1:A:1035:ILE:HG12	1:A:1049:MET:HB3	2.04	0.40
1:A:1053:ASP:OD1	1:A:1053:ASP:N	2.50	0.40
1:A:1068:GLY:C	3:A:3141:HOH:O	2.63	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	296/356 (83%)	280 (95%)	13 (4%)	3 (1%)	12 5

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1058	SER
1	A	916	GLU
1	A	939	LEU

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	250/308 (81%)	229 (92%)	21 (8%)	10 4

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	855	LEU
1	A	878	GLU
1	A	909	LEU
1	A	920	LEU
1	A	939	LEU
1	A	962	LYS
1	A	977	LEU
1	A	985	SER
1	A	1006	LEU
1	A	1031	ILE
1	A	1046	LEU
1	A	1049	MET
1	A	1059	VAL
1	A	1073	VAL
1	A	1074	LEU
1	A	1120	GLU
1	A	1123	VAL
1	A	1144	LEU
1	A	1163	LYS
1	A	1171	LEU
1	A	1176	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	913	HIS
1	A	1004	GLN
1	A	1096	ASN
1	A	1114	GLN
1	A	1158	HIS
1	A	1175	GLN

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](#)

5 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	$\# Z > 2$
2	GOL	A	2188	-	5,5,5	0.52	0	5,5,5	1.00	0
2	GOL	A	2189	-	5,5,5	0.31	0	5,5,5	0.42	0
2	GOL	A	2190	-	5,5,5	0.48	0	5,5,5	0.57	0
2	GOL	A	2187	-	5,5,5	0.40	0	5,5,5	0.74	0
2	GOL	A	2191	-	5,5,5	0.35	0	5,5,5	0.51	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	2188	-	-	3/4/4/4	-
2	GOL	A	2189	-	-	4/4/4/4	-
2	GOL	A	2190	-	-	2/4/4/4	-
2	GOL	A	2187	-	-	0/4/4/4	-
2	GOL	A	2191	-	-	0/4/4/4	-

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
2	A	2189	GOL	O1-C1-C2-C3
2	A	2190	GOL	O1-C1-C2-C3
2	A	2189	GOL	O1-C1-C2-O2
2	A	2188	GOL	O1-C1-C2-C3
2	A	2189	GOL	C1-C2-C3-O3
2	A	2188	GOL	O1-C1-C2-O2
2	A	2189	GOL	O2-C2-C3-O3
2	A	2190	GOL	O1-C1-C2-O2
2	A	2188	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2188	GOL	2	0
2	A	2190	GOL	2	0
2	A	2191	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

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	306/356 (85%)	0.52	25 (8%) 17 18	15, 35, 56, 76	3 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	949	CYS	5.4
1	A	1103	VAL	5.0
1	A	1115	ALA	3.5
1	A	1112	PRO	3.4
1	A	1116	GLY	3.3
1	A	1102	SER	3.1
1	A	1113	GLY	3.1
1	A	854	ASN	2.8
1	A	1148	THR	2.8
1	A	1064	TYR	2.7
1	A	1088	PRO	2.7
1	A	1101	LEU	2.7
1	A	1057	ALA	2.6
1	A	985	SER	2.6
1	A	1109	CYS	2.5
1	A	984	LEU	2.5
1	A	1075	SER	2.3
1	A	1105	VAL	2.3
1	A	868	CYS	2.3
1	A	1144	LEU	2.3
1	A	1104	GLY	2.3
1	A	1110	LEU	2.3
1	A	956	GLU	2.0
1	A	1094	VAL	2.0
1	A	948	PHE	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	A	2189	6/6	0.80	0.12	37,47,48,48	0
2	GOL	A	2190	6/6	0.87	0.11	36,41,43,45	0
2	GOL	A	2187	6/6	0.88	0.09	37,39,43,47	0
2	GOL	A	2191	6/6	0.89	0.11	42,51,55,58	0
2	GOL	A	2188	6/6	0.92	0.12	23,35,42,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.