



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2022 – 01:10 pm BST

PDB ID : 7PLN
Title : Structure of the APCbeta domain of Plasmodium vivax perforin-like protein 1
Authors : Williams, S.I.; Ni, T.; Yu, X.; Gilbert, R.J.C.
Deposited on : 2021-08-31
Resolution : 3.15 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
Xtriage (Phenix) : 1.13
EDS : 2.29
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.29

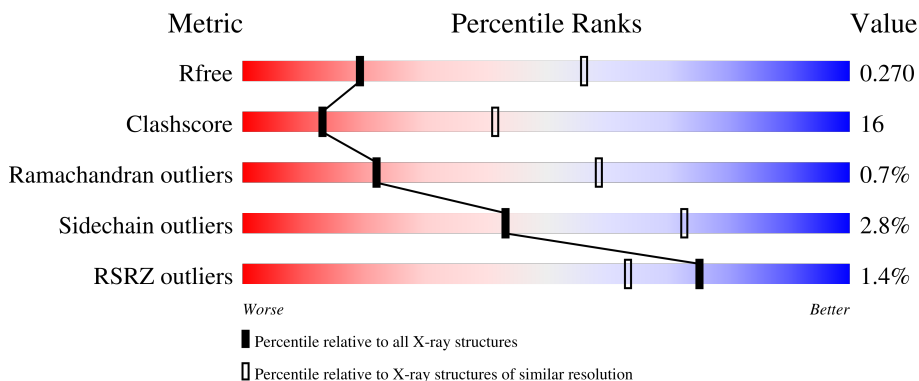
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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}	130704	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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	271	
1	B	271	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3682 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 Sporozoite micronemal protein essential for cell traversal, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	242	Total	C	N	O	S	0	0	0
			1846	1168	299	362	17			
1	B	241	Total	C	N	O	S	0	0	0
			1836	1160	300	359	17			

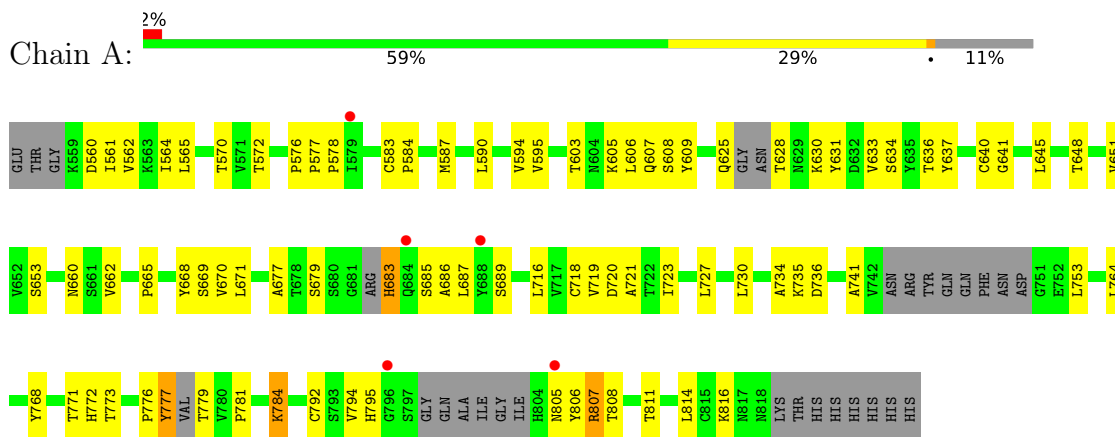
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	556	GLU	-	expression tag	UNP A0A1G4H7E1
A	557	THR	-	expression tag	UNP A0A1G4H7E1
A	558	GLY	-	expression tag	UNP A0A1G4H7E1
A	819	LYS	-	expression tag	UNP A0A1G4H7E1
A	820	THR	-	expression tag	UNP A0A1G4H7E1
A	821	HIS	-	expression tag	UNP A0A1G4H7E1
A	822	HIS	-	expression tag	UNP A0A1G4H7E1
A	823	HIS	-	expression tag	UNP A0A1G4H7E1
A	824	HIS	-	expression tag	UNP A0A1G4H7E1
A	825	HIS	-	expression tag	UNP A0A1G4H7E1
A	826	HIS	-	expression tag	UNP A0A1G4H7E1
B	556	GLU	-	expression tag	UNP A0A1G4H7E1
B	557	THR	-	expression tag	UNP A0A1G4H7E1
B	558	GLY	-	expression tag	UNP A0A1G4H7E1
B	819	LYS	-	expression tag	UNP A0A1G4H7E1
B	820	THR	-	expression tag	UNP A0A1G4H7E1
B	821	HIS	-	expression tag	UNP A0A1G4H7E1
B	822	HIS	-	expression tag	UNP A0A1G4H7E1
B	823	HIS	-	expression tag	UNP A0A1G4H7E1
B	824	HIS	-	expression tag	UNP A0A1G4H7E1
B	825	HIS	-	expression tag	UNP A0A1G4H7E1
B	826	HIS	-	expression tag	UNP A0A1G4H7E1

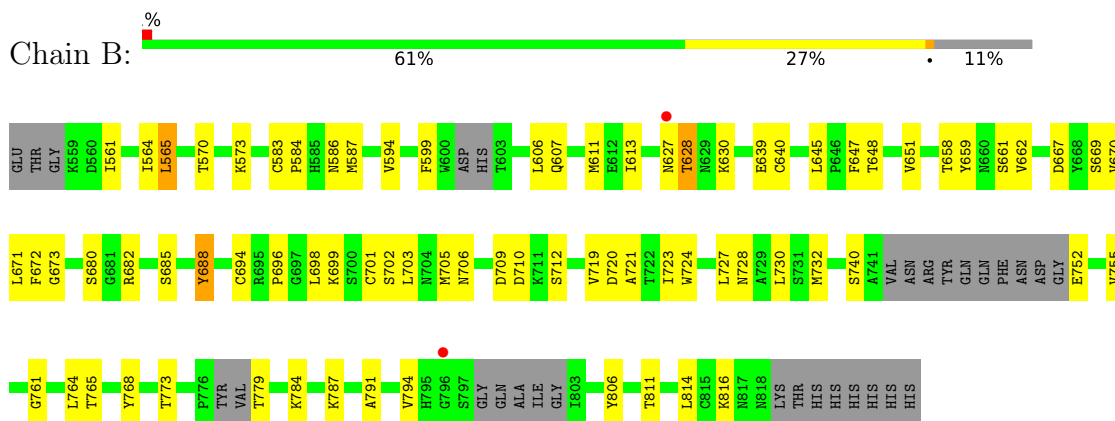
3 Residue-property plots [i](#)

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: Sporozoite micronemal protein essential for cell traversal, putative



- Molecule 1: Sporozoite micronemal protein essential for cell traversal, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.74Å 42.02Å 87.22Å 90.00° 104.95° 90.00°	Depositor
Resolution (Å)	63.01 – 3.15 63.01 – 3.15	Depositor EDS
% Data completeness (in resolution range)	98.4 (63.01-3.15) 98.4 (63.01-3.15)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 3.13Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.212 , 0.269 0.211 , 0.270	Depositor DCC
R_{free} test set	431 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	85.9	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3682	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.38% 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](#)

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	0.61	1/1887 (0.1%)	0.77	1/2561 (0.0%)
1	B	0.54	1/1876 (0.1%)	0.78	1/2546 (0.0%)
All	All	0.57	2/3763 (0.1%)	0.77	2/5107 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	583	CYS	CB-SG	-5.47	1.72	1.81
1	A	583	CYS	CB-SG	-5.40	1.73	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	583	CYS	CA-CB-SG	-5.11	104.80	114.00
1	B	565	LEU	CB-CG-CD2	-5.02	102.46	111.00

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	1846	0	1773	62	1
1	B	1836	0	1775	58	1
All	All	3682	0	3548	118	1

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 16.

All (118) 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:794:VAL:HG12	1:A:795:HIS:H	1.42	0.84
1:A:677:ALA:O	1:A:689:SER:HA	1.79	0.82
1:A:625:GLN:O	1:A:628:THR:N	2.18	0.76
1:B:768:TYR:HB2	1:B:794:VAL:HG22	1.69	0.75
1:A:645:LEU:HB2	1:A:648:THR:HG21	1.69	0.73
1:A:603:THR:HB	1:A:605:LYS:HD3	1.71	0.71
1:A:670:VAL:HG22	1:A:716:LEU:HD13	1.76	0.68
1:A:773:THR:OG1	1:A:805:ASN:HB3	1.93	0.68
1:A:687:LEU:HD13	1:A:807:ARG:HA	1.75	0.67
1:B:761:GLY:HA2	1:B:816:LYS:O	1.94	0.67
1:A:561:ILE:HG12	1:A:671:LEU:HD11	1.77	0.66
1:A:687:LEU:HD23	1:A:689:SER:OG	1.95	0.66
1:A:792:CYS:HB3	1:A:811:THR:HG21	1.76	0.65
1:B:673:GLY:H	1:B:732:MET:HE1	1.61	0.65
1:A:768:TYR:HB2	1:A:794:VAL:HG22	1.78	0.64
1:A:771:THR:O	1:A:807:ARG:HB2	1.97	0.64
1:B:669:SER:HB2	1:B:721:ALA:HB2	1.80	0.63
1:A:631:TYR:HE2	1:A:634:SER:HB3	1.65	0.62
1:A:730:LEU:HD23	1:A:814:LEU:HD21	1.81	0.62
1:A:687:LEU:CD1	1:A:807:ARG:HA	2.31	0.61
1:B:651:VAL:HG11	1:B:662:VAL:HG21	1.84	0.60
1:B:570:THR:HG23	1:B:787:LYS:NZ	2.18	0.58
1:B:740:SER:HA	1:B:806:TYR:CE1	2.38	0.58
1:B:561:ILE:HB	1:B:724:TRP:HE3	1.67	0.58
1:B:658:THR:HA	1:B:705:MET:HE2	1.86	0.57
1:B:564:ILE:CD1	1:B:645:LEU:HD21	2.35	0.57
1:A:735:LYS:HG2	1:A:753:LEU:HD11	1.87	0.56
1:A:679:SER:OG	1:A:679:SER:O	2.21	0.56
1:B:565:LEU:O	1:B:787:LYS:HE2	2.06	0.55
1:A:565:LEU:CB	1:A:764:LEU:HD12	2.36	0.54
1:B:627:ASN:O	1:B:628:THR:OG1	2.22	0.54
1:B:794:VAL:HG23	1:B:811:THR:HG23	1.89	0.54
1:B:688:TYR:OH	1:B:709:ASP:OD1	2.16	0.54
1:B:648:THR:HG22	1:B:719:VAL:HG12	1.90	0.53
1:B:673:GLY:H	1:B:732:MET:CE	2.22	0.53
1:A:584:PRO:HB2	1:A:587:MET:HG3	1.89	0.53
1:A:776:PRO:HG2	1:A:779:THR:HG21	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:565:LEU:HB3	1:A:764:LEU:HD12	1.91	0.53
1:A:561:ILE:HD12	1:A:564:ILE:HD11	1.91	0.52
1:B:639:GLU:OE1	1:B:765:THR:OG1	2.18	0.52
1:A:590:LEU:HD11	1:A:641:GLY:N	2.26	0.51
1:B:647:PHE:HB3	1:B:720:ASP:HB2	1.92	0.51
1:B:606:LEU:HG	1:B:607:GLN:N	2.26	0.51
1:B:671:LEU:HD13	1:B:719:VAL:HG11	1.92	0.51
1:A:794:VAL:HG12	1:A:795:HIS:N	2.19	0.50
1:B:570:THR:HG22	1:B:639:GLU:OE1	2.11	0.50
1:A:741:ALA:HB2	1:A:806:TYR:HB3	1.94	0.50
1:B:720:ASP:HB3	1:B:723:ILE:HD13	1.92	0.50
1:B:606:LEU:HD22	1:B:773:THR:HG22	1.93	0.50
1:B:659:TYR:CZ	1:B:661:SER:HB2	2.45	0.50
1:A:719:VAL:CG2	1:A:727:LEU:HD11	2.42	0.50
1:A:653:SER:OG	1:A:662:VAL:HG22	2.13	0.49
1:B:594:VAL:O	1:B:611:MET:HA	2.13	0.48
1:A:572:THR:HG23	1:A:784:LYS:HB3	1.95	0.48
1:B:703:LEU:HD22	1:B:712:SER:OG	2.14	0.48
1:B:671:LEU:HD23	1:B:672:PHE:HB2	1.95	0.48
1:A:630:LYS:HG2	1:A:631:TYR:N	2.29	0.47
1:A:687:LEU:HD22	1:A:808:THR:HG23	1.97	0.47
1:B:673:GLY:HA3	1:B:694:CYS:HB2	1.97	0.47
1:A:590:LEU:H	1:A:590:LEU:HD12	1.79	0.46
1:A:651:VAL:HG11	1:A:662:VAL:HG11	1.96	0.46
1:A:730:LEU:HB3	1:A:814:LEU:HD11	1.97	0.46
1:B:561:ILE:HB	1:B:724:TRP:CE3	2.49	0.46
1:B:570:THR:OG1	1:B:784:LYS:HB2	2.15	0.46
1:B:701:CYS:SG	1:B:702:SER:N	2.89	0.46
1:B:740:SER:HA	1:B:806:TYR:CD1	2.50	0.46
1:B:630:LYS:O	1:B:630:LYS:HG3	2.15	0.46
1:B:613:ILE:O	1:B:613:ILE:HG13	2.16	0.46
1:B:727:LEU:O	1:B:730:LEU:HB2	2.17	0.45
1:B:570:THR:HG23	1:B:787:LYS:HZ1	1.81	0.45
1:A:764:LEU:HD21	1:A:816:LYS:HB3	1.97	0.45
1:B:680:SER:H	1:B:710:ASP:CG	2.20	0.45
1:B:728:ASN:C	1:B:730:LEU:H	2.19	0.44
1:B:688:TYR:C	1:B:688:TYR:CD1	2.91	0.44
1:A:719:VAL:HG23	1:A:727:LEU:HD11	1.99	0.44
1:A:669:SER:HB2	1:A:721:ALA:HB2	1.99	0.44
1:A:779:THR:HG23	1:A:779:THR:O	2.18	0.44
1:A:683:HIS:CE1	1:A:685:SER:HB2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:670:VAL:N	1:B:698:LEU:O	2.41	0.43
1:B:706:ASN:HB2	1:B:709:ASP:OD2	2.18	0.43
1:A:562:VAL:HA	1:A:764:LEU:HD11	2.01	0.43
1:B:599:PHE:HD1	1:B:606:LEU:HA	1.83	0.43
1:A:576:PRO:HA	1:A:633:VAL:HA	2.01	0.43
1:A:687:LEU:HD12	1:A:773:THR:HG23	1.99	0.43
1:B:584:PRO:O	1:B:587:MET:HB2	2.18	0.43
1:A:603:THR:CB	1:A:605:LYS:HD3	2.44	0.43
1:A:565:LEU:HB2	1:A:764:LEU:HD12	1.99	0.43
1:A:686:ALA:HB3	1:A:773:THR:HG21	2.01	0.43
1:B:667:ASP:OD2	1:B:699:LYS:NZ	2.49	0.43
1:B:727:LEU:HD12	1:B:727:LEU:HA	1.59	0.43
1:B:732:MET:SD	1:B:814:LEU:HD13	2.59	0.43
1:A:794:VAL:CG1	1:A:795:HIS:H	2.24	0.42
1:A:570:THR:HG23	1:A:637:TYR:CE2	2.55	0.42
1:B:755:VAL:O	1:B:791:ALA:HB1	2.20	0.42
1:A:687:LEU:HD23	1:A:689:SER:CB	2.48	0.42
1:B:630:LYS:HB2	1:B:630:LYS:HE3	1.41	0.42
1:A:577:PRO:HA	1:A:578:PRO:HA	1.89	0.42
1:A:687:LEU:HG	1:A:773:THR:HG22	2.01	0.42
1:A:606:LEU:HD21	1:A:609:TYR:CE2	2.55	0.42
1:A:671:LEU:C	1:A:671:LEU:HD23	2.40	0.42
1:B:564:ILE:HD12	1:B:645:LEU:HD21	2.02	0.42
1:B:740:SER:HA	1:B:806:TYR:HE1	1.84	0.42
1:B:645:LEU:HB3	1:B:648:THR:HG21	2.02	0.41
1:A:594:VAL:HG22	1:A:636:THR:HG23	2.01	0.41
1:A:665:PRO:HD2	1:A:718:CYS:SG	2.60	0.41
1:A:595:VAL:O	1:A:634:SER:HA	2.20	0.41
1:A:687:LEU:HD11	1:A:772:HIS:C	2.40	0.41
1:B:814:LEU:HD12	1:B:814:LEU:HA	1.82	0.41
1:B:627:ASN:OD1	1:B:627:ASN:N	2.53	0.41
1:B:764:LEU:O	1:B:787:LYS:HE3	2.20	0.41
1:A:734:ALA:HA	1:A:811:THR:O	2.21	0.41
1:B:670:VAL:O	1:B:696:PRO:HA	2.19	0.41
1:A:736:ASP:H	1:B:685:SER:HB3	1.86	0.41
1:A:668:TYR:CE1	1:A:720:ASP:HB2	2.56	0.40
1:B:573:LYS:HE3	1:B:573:LYS:HB2	1.73	0.40
1:A:576:PRO:HG3	1:A:777:TYR:HD2	1.86	0.40
1:A:607:GLN:HG2	1:A:608:SER:N	2.36	0.40
1:A:735:LYS:HG3	1:B:685:SER:HB3	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:776:PRO:O	1:B:779:THR:OG1[2_554]	2.19	0.01

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	230/271 (85%)	198 (86%)	30 (13%)	2 (1%)	17	53
1	B	231/271 (85%)	199 (86%)	31 (13%)	1 (0%)	34	68
All	All	461/542 (85%)	397 (86%)	61 (13%)	3 (1%)	22	59

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	781	PRO
1	A	723	ILE
1	B	628	THR

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	213/237 (90%)	206 (97%)	7 (3%)	38	69
1	B	212/237 (90%)	207 (98%)	5 (2%)	49	76
All	All	425/474 (90%)	413 (97%)	12 (3%)	43	73

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	560	ASP
1	A	640	CYS
1	A	660	ASN
1	A	683	HIS
1	A	777	TYR
1	A	784	LYS
1	A	807	ARG
1	B	586	ASN
1	B	640	CYS
1	B	682	ARG
1	B	688	TYR
1	B	752	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

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	242/271 (89%)	0.06	5 (2%) 63 49	45, 82, 129, 160	0
1	B	241/271 (88%)	-0.05	2 (0%) 86 78	51, 78, 131, 187	0
All	All	483/542 (89%)	0.01	7 (1%) 75 63	45, 80, 130, 187	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	688	TYR	3.0
1	A	805	ASN	2.8
1	B	796	GLY	2.5
1	A	684	GLN	2.5
1	B	627	ASN	2.4
1	A	796	GLY	2.1
1	A	579	ILE	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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.