



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:53 PM UTC

PDB ID : 7LGN / pdb_00007lgn
Title : Cyanophycin synthetase 1 from *T. morbirosei*
Authors : Sharon, I.; Schmeing, T.M.
Deposited on : 2021-01-20
Resolution : 3.10 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
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

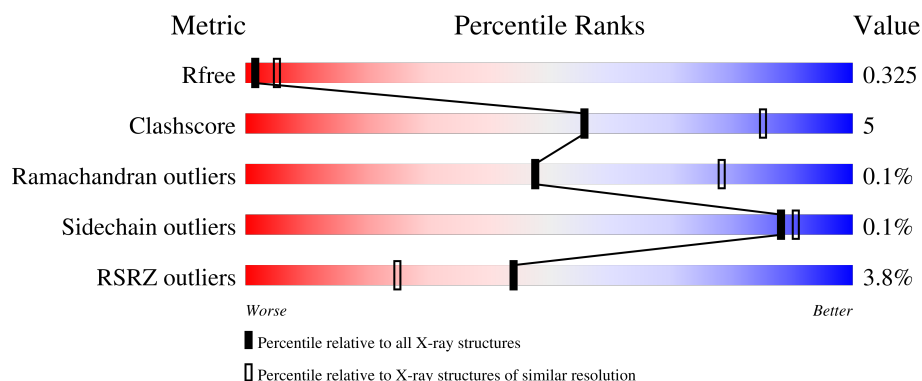
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	909	
1	B	909	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11790 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 Cyanophycin synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	706	Total	C	N	O	S	0	0	0
			5321	3322	945	1025	29			
1	B	850	Total	C	N	O	S	0	0	0
			6469	4040	1154	1236	39			

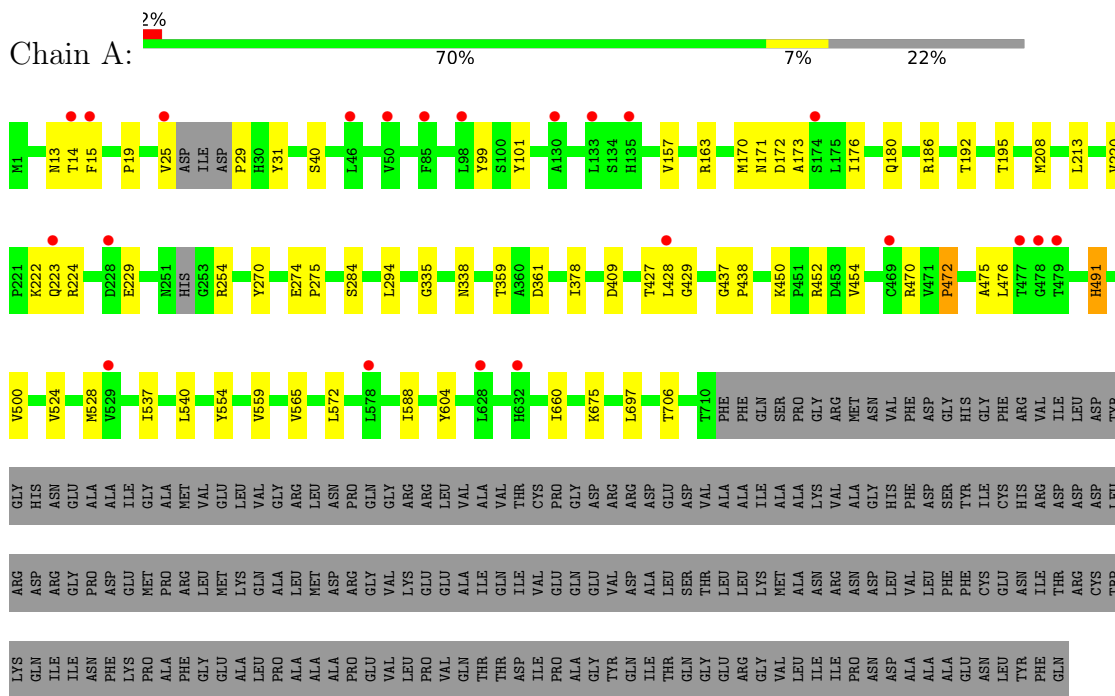
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	901	ALA	-	expression tag	UNP A0A095T5Z8
A	902	ALA	-	expression tag	UNP A0A095T5Z8
A	903	ALA	-	expression tag	UNP A0A095T5Z8
A	904	GLU	-	expression tag	UNP A0A095T5Z8
A	905	ASN	-	expression tag	UNP A0A095T5Z8
A	906	LEU	-	expression tag	UNP A0A095T5Z8
A	907	TYR	-	expression tag	UNP A0A095T5Z8
A	908	PHE	-	expression tag	UNP A0A095T5Z8
A	909	GLN	-	expression tag	UNP A0A095T5Z8
B	901	ALA	-	expression tag	UNP A0A095T5Z8
B	902	ALA	-	expression tag	UNP A0A095T5Z8
B	903	ALA	-	expression tag	UNP A0A095T5Z8
B	904	GLU	-	expression tag	UNP A0A095T5Z8
B	905	ASN	-	expression tag	UNP A0A095T5Z8
B	906	LEU	-	expression tag	UNP A0A095T5Z8
B	907	TYR	-	expression tag	UNP A0A095T5Z8
B	908	PHE	-	expression tag	UNP A0A095T5Z8
B	909	GLN	-	expression tag	UNP A0A095T5Z8

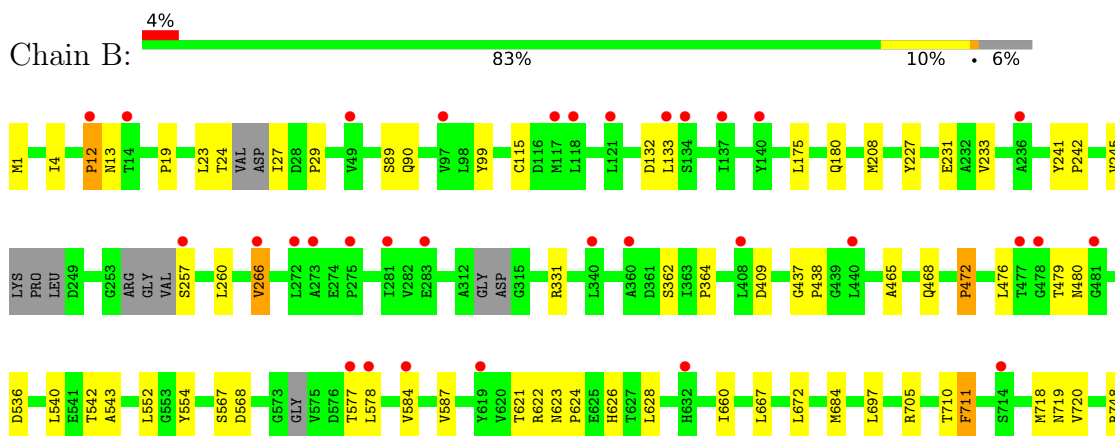
3 Residue-property plots

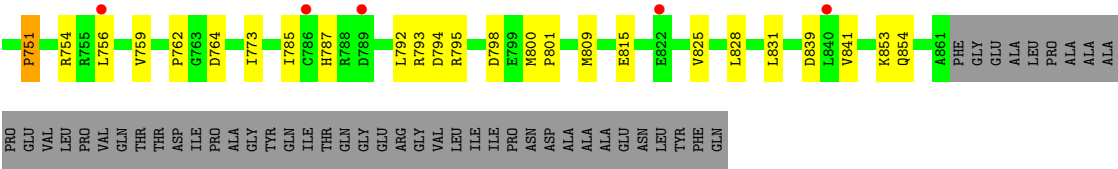
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Cyanophycin synthase



• Molecule 1: Cyanophycin synthase





4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	222.56Å 222.56Å 105.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	72.31 – 3.10 72.31 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (72.31-3.10) 94.1 (72.31-3.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.33 (at 3.07Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.270 , 0.326 0.274 , 0.325	Depositor DCC
R_{free} test set	2413 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	79.5	Xtriage
Anisotropy	0.798	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 89.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.015 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11790	wwPDB-VP
Average B, all atoms (Å ²)	130.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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

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	0.60	0/5399	1.17	7/7324 (0.1%)
1	B	0.57	0/6566	1.17	13/8895 (0.1%)
All	All	0.58	0/11965	1.17	20/16219 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	792	LEU	N-CA-C	15.25	130.92	111.24
1	A	572	LEU	N-CA-C	-7.54	104.58	114.31
1	B	792	LEU	N-CA-CB	-7.44	99.13	110.07
1	A	378	ILE	N-CA-C	-7.29	105.54	111.81
1	B	711	PHE	CB-CA-C	-7.24	101.45	111.73
1	B	364	PRO	N-CA-C	6.79	118.99	110.70
1	A	675	LYS	N-CA-C	-6.12	102.96	110.61
1	B	719	ASN	CA-CB-CG	-6.10	106.50	112.60
1	A	491	HIS	CA-CB-CG	-5.64	108.16	113.80
1	B	12	PRO	N-CA-C	5.51	123.83	112.47
1	B	362	SER	N-CA-C	5.49	117.44	108.55
1	B	133	LEU	N-CA-C	-5.39	106.54	113.23
1	B	331	ARG	N-CA-C	-5.38	106.89	113.50
1	B	795	ARG	N-CA-C	-5.33	104.20	111.56
1	B	751	PRO	CB-CA-C	-5.33	104.59	111.46
1	A	40	SER	N-CA-C	-5.25	106.82	114.12
1	A	472	PRO	N-CA-C	-5.23	103.62	111.57
1	A	224	ARG	CB-CA-C	-5.17	104.93	111.86
1	B	227	TYR	CB-CA-C	5.11	118.76	110.22
1	B	472	PRO	N-CA-C	-5.09	103.84	111.57

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	5321	0	5386	45	0
1	B	6469	0	6502	74	0
All	All	11790	0	11888	112	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 (112) 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:B:759:VAL:HG21	1:B:831:LEU:HD11	1.64	0.79
1:B:12:PRO:HD2	1:B:552:LEU:HB2	1.68	0.75
1:B:759:VAL:CG2	1:B:831:LEU:HD11	2.18	0.73
1:A:25:VAL:HB	1:A:31:TYR:HD1	1.55	0.72
1:B:825:VAL:HG13	1:B:854:GLN:NE2	2.04	0.71
1:B:4:ILE:HD11	1:B:24:THR:CG2	2.22	0.70
1:A:171:ASN:HB2	1:B:208:MET:CE	2.23	0.69
1:B:4:ILE:HD11	1:B:24:THR:HG23	1.73	0.69
1:B:825:VAL:HG13	1:B:854:GLN:HE22	1.57	0.69
1:A:180:GLN:HB2	1:A:472:PRO:HD2	1.74	0.68
1:B:809:MET:SD	1:B:815:GLU:HB2	2.35	0.66
1:B:828:LEU:HD23	1:B:854:GLN:OE1	1.95	0.65
1:B:793:ARG:HG2	1:B:794:ASP:N	2.12	0.64
1:B:754:ARG:HB2	1:B:839:ASP:OD1	1.98	0.63
1:A:500:VAL:HG23	1:A:537:ILE:HG23	1.81	0.63
1:B:793:ARG:HG2	1:B:794:ASP:H	1.63	0.63
1:A:660:ILE:CG2	1:A:697:LEU:HD11	2.31	0.60
1:B:245:VAL:O	1:B:257:SER:OG	2.20	0.60
1:A:19:PRO:HD2	1:A:99:TYR:O	2.02	0.59
1:A:192:THR:O	1:A:195:THR:HG22	2.04	0.56
1:B:584:VAL:O	1:B:587:VAL:HG22	2.05	0.56
1:A:170:MET:HE1	1:A:186:ARG:HG2	1.87	0.55
1:B:759:VAL:HG22	1:B:785:ILE:HB	1.88	0.55
1:B:622:ARG:C	1:B:624:PRO:HD2	2.31	0.55
1:B:536:ASP:O	1:B:536:ASP:OD1	2.25	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:PRO:C	1:A:31:TYR:H	2.15	0.54
1:B:180:GLN:HB2	1:B:472:PRO:CD	2.38	0.54
1:A:180:GLN:HB2	1:A:472:PRO:CD	2.38	0.53
1:A:19:PRO:HD3	1:A:101:TYR:CD1	2.43	0.53
1:A:284:SER:HB3	1:B:853:LYS:HZ3	1.74	0.53
1:B:465:ALA:HA	1:B:468:GLN:HB2	1.90	0.53
1:B:754:ARG:CB	1:B:839:ASP:OD1	2.55	0.52
1:B:1:MET:H3	1:B:115:CYS:HB3	1.74	0.52
1:B:132:ASP:O	1:B:132:ASP:OD1	2.28	0.51
1:A:565:VAL:HG21	1:A:604:TYR:HB2	1.92	0.51
1:B:798:ASP:OD1	1:B:798:ASP:N	2.43	0.51
1:B:762:PRO:HG3	1:B:773:ILE:HD13	1.93	0.50
1:A:491:HIS:NE2	1:A:706:THR:O	2.46	0.49
1:B:476:LEU:CD1	1:B:540:LEU:O	2.61	0.49
1:B:720:VAL:HG23	1:B:720:VAL:O	2.12	0.49
1:A:14:THR:OG1	1:A:15:PHE:N	2.46	0.49
1:B:89:SER:OG	1:B:90:GLN:N	2.45	0.48
1:B:621:THR:O	1:B:684:MET:HE1	2.13	0.48
1:A:427:THR:O	1:A:429:GLY:N	2.47	0.48
1:A:524:VAL:O	1:A:528:MET:HG2	2.13	0.48
1:A:213:LEU:HB2	1:A:220:VAL:HG21	1.95	0.48
1:A:171:ASN:HB2	1:B:208:MET:HE3	1.95	0.47
1:B:759:VAL:HG12	1:B:787:HIS:ND1	2.29	0.47
1:B:180:GLN:HB2	1:B:472:PRO:HD2	1.95	0.47
1:B:479:THR:OG1	1:B:480:ASN:N	2.47	0.47
1:B:764:ASP:OD1	1:B:764:ASP:N	2.36	0.47
1:B:19:PRO:HD2	1:B:99:TYR:O	2.15	0.47
1:A:428:LEU:HD22	1:B:853:LYS:HE2	1.96	0.47
1:B:23:LEU:C	1:B:23:LEU:HD12	2.40	0.46
1:A:359:THR:HG23	1:A:361:ASP:H	1.80	0.46
1:A:284:SER:HB3	1:B:853:LYS:NZ	2.30	0.46
1:A:294:LEU:HD21	1:A:454:VAL:HG11	1.98	0.46
1:B:4:ILE:CD1	1:B:24:THR:HG23	2.45	0.46
1:B:437:GLY:N	1:B:438:PRO:CD	2.78	0.46
1:B:542:THR:OG1	1:B:543:ALA:N	2.49	0.46
1:A:476:LEU:HD23	1:A:588:ILE:HG12	1.99	0.45
1:B:180:GLN:HB2	1:B:472:PRO:HD3	1.98	0.45
1:B:762:PRO:HG3	1:B:773:ILE:CD1	2.46	0.45
1:B:476:LEU:HD13	1:B:540:LEU:O	2.16	0.45
1:A:476:LEU:HD13	1:A:540:LEU:O	2.17	0.45
1:B:793:ARG:CG	1:B:794:ASP:H	2.29	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:831:LEU:HD13	1:B:841:VAL:HG13	1.99	0.45
1:A:437:GLY:N	1:A:438:PRO:CD	2.80	0.45
1:B:787:HIS:CG	1:B:787:HIS:O	2.70	0.45
1:B:751:PRO:HG2	1:B:756:LEU:HD21	1.99	0.44
1:A:13:ASN:N	1:A:13:ASN:OD1	2.50	0.44
1:A:660:ILE:HG22	1:A:697:LEU:HD11	1.98	0.44
1:B:241:TYR:N	1:B:242:PRO:CD	2.80	0.44
1:B:12:PRO:HB2	1:B:552:LEU:HD13	1.99	0.44
1:B:577:THR:OG1	1:B:578:LEU:N	2.51	0.44
1:B:787:HIS:O	1:B:787:HIS:CD2	2.70	0.44
1:B:626:HIS:CE1	1:B:628:LEU:HB3	2.52	0.43
1:B:793:ARG:CG	1:B:794:ASP:N	2.79	0.43
1:A:163:ARG:O	1:A:470:ARG:NH1	2.52	0.43
1:A:172:ASP:CG	1:A:173:ALA:H	2.26	0.43
1:B:409:ASP:OD1	1:B:409:ASP:N	2.52	0.43
1:B:27:ILE:HG22	1:B:29:PRO:HD2	2.00	0.43
1:B:710:THR:O	1:B:711:PHE:HB2	2.19	0.43
1:A:450:LYS:HB2	1:A:452:ARG:HH12	1.83	0.43
1:B:260:LEU:HB3	1:B:266:VAL:HG22	2.01	0.43
1:B:623:ASN:N	1:B:624:PRO:HD2	2.34	0.42
1:A:195:THR:O	1:A:195:THR:HG23	2.19	0.42
1:B:567:SER:OG	1:B:568:ASP:N	2.51	0.42
1:A:335:GLY:H	1:A:338:ASN:HD22	1.66	0.42
1:B:23:LEU:HD12	1:B:24:THR:N	2.35	0.42
1:B:718:MET:HG2	1:B:718:MET:O	2.20	0.42
1:A:222:LYS:O	1:A:223:GLN:C	2.62	0.41
1:B:672:LEU:CD1	1:B:748:ARG:NH2	2.83	0.41
1:B:231:GLU:O	1:B:233:VAL:N	2.53	0.41
1:B:554:TYR:CD1	1:B:554:TYR:C	2.98	0.41
1:A:409:ASP:N	1:A:409:ASP:OD1	2.53	0.41
1:B:476:LEU:HD22	1:B:554:TYR:OH	2.20	0.41
1:A:157:VAL:CG2	1:A:176:ILE:HD13	2.50	0.41
1:A:171:ASN:HB2	1:B:208:MET:SD	2.60	0.41
1:A:475:ALA:HA	1:A:559:VAL:O	2.21	0.41
1:B:660:ILE:HG21	1:B:697:LEU:HD11	2.01	0.41
1:B:241:TYR:N	1:B:242:PRO:HD3	2.36	0.41
1:B:623:ASN:N	1:B:624:PRO:CD	2.82	0.41
1:B:800:MET:HB2	1:B:801:PRO:HD3	2.03	0.41
1:A:29:PRO:C	1:A:31:TYR:N	2.79	0.41
1:A:554:TYR:CD1	1:A:554:TYR:C	2.99	0.41
1:A:208:MET:HE1	1:B:175:LEU:HD23	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:GLU:HB3	1:A:270:TYR:OH	2.21	0.41
1:A:274:GLU:N	1:A:275:PRO:HD2	2.36	0.41
1:A:254:ARG:O	1:A:254:ARG:HG2	2.21	0.40
1:A:213:LEU:CB	1:A:220:VAL:HG21	2.51	0.40
1:B:667:LEU:HD13	1:B:705:ARG:CZ	2.51	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	700/909 (77%)	670 (96%)	30 (4%)	0	100	100
1	B	838/909 (92%)	790 (94%)	47 (6%)	1 (0%)	48	78
All	All	1538/1818 (85%)	1460 (95%)	77 (5%)	1 (0%)	48	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	13	ASN

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	569/735 (77%)	569 (100%)	0	100	100
1	B	690/735 (94%)	689 (100%)	1 (0%)	88	90
All	All	1259/1470 (86%)	1258 (100%)	1 (0%)	88	90

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	266	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	435	ASN
1	A	443	HIS
1	B	13	ASN
1	B	177	GLN
1	B	348	GLN
1	B	443	HIS
1	B	611	HIS
1	B	806	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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

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	706/909 (77%)	0.23	22 (3%)	51 30	48, 102, 205, 250	0
1	B	850/909 (93%)	0.43	37 (4%)	39 21	58, 132, 237, 273	0
All	All	1556/1818 (85%)	0.34	59 (3%)	44 25	48, 115, 227, 273	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	118	LEU	4.8
1	A	478	GLY	3.9
1	B	117	MET	3.7
1	B	49	VAL	3.7
1	B	281	ILE	3.7
1	A	628	LEU	3.7
1	B	121	LEU	3.6
1	B	789	ASP	3.6
1	A	632	HIS	3.5
1	B	283	GLU	3.4
1	B	756	LEU	3.4
1	B	133	LEU	3.3
1	B	266	VAL	3.3
1	A	428	LEU	3.2
1	A	477	THR	3.0
1	B	273	ALA	2.9
1	A	174	SER	2.9
1	B	272	LEU	2.9
1	B	632	HIS	2.8
1	B	275	PRO	2.8
1	B	440	LEU	2.8
1	A	15	PHE	2.8
1	B	481	GLY	2.7
1	B	14	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	137	ILE	2.6
1	B	577	THR	2.6
1	A	14	THR	2.5
1	A	46	LEU	2.5
1	A	130	ALA	2.5
1	A	98	LEU	2.4
1	A	578	LEU	2.4
1	A	25	VAL	2.4
1	B	478	GLY	2.4
1	B	97	VAL	2.3
1	B	236	ALA	2.3
1	B	619	TYR	2.3
1	B	477	THR	2.3
1	B	578	LEU	2.3
1	A	223	GLN	2.3
1	B	584	VAL	2.3
1	A	469	CYS	2.2
1	B	786	CYS	2.2
1	A	85	PHE	2.2
1	A	529	VAL	2.2
1	B	134	SER	2.2
1	A	228	ASP	2.2
1	A	135	HIS	2.2
1	B	340	LEU	2.1
1	A	479	THR	2.1
1	B	840	LEU	2.1
1	B	360	ALA	2.1
1	B	822	GLU	2.1
1	B	714	SER	2.1
1	B	140	TYR	2.0
1	A	50	VAL	2.0
1	A	133	LEU	2.0
1	B	408	LEU	2.0
1	B	257	SER	2.0
1	B	12	PRO	2.0

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

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.