



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 24, 2026 – 09:22 PM EDT

PDB ID : 1BJT / pdb_00001bjt
Title : TOPOISOMERASE II RESIDUES 409-1201
Authors : Fass, D.; Bogden, C.E.; Berger, J.M.
Deposited on : 1998-06-29
Resolution : 2.50 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

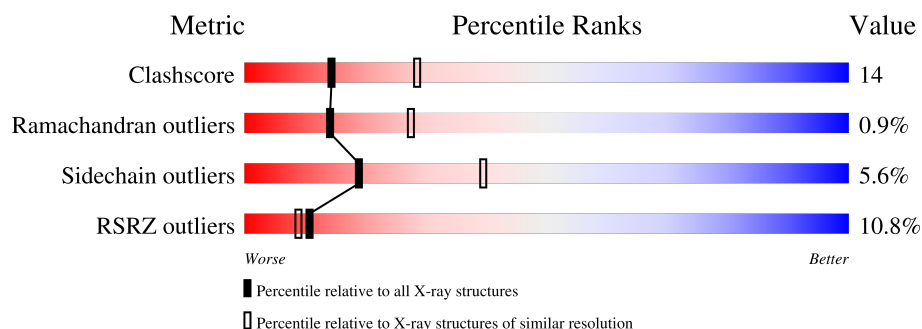
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	793	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5815 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 TOPOISOMERASE II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	679	5598	3621	939	1018	20	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	547	LEU	PRO	conflict	UNP P06786

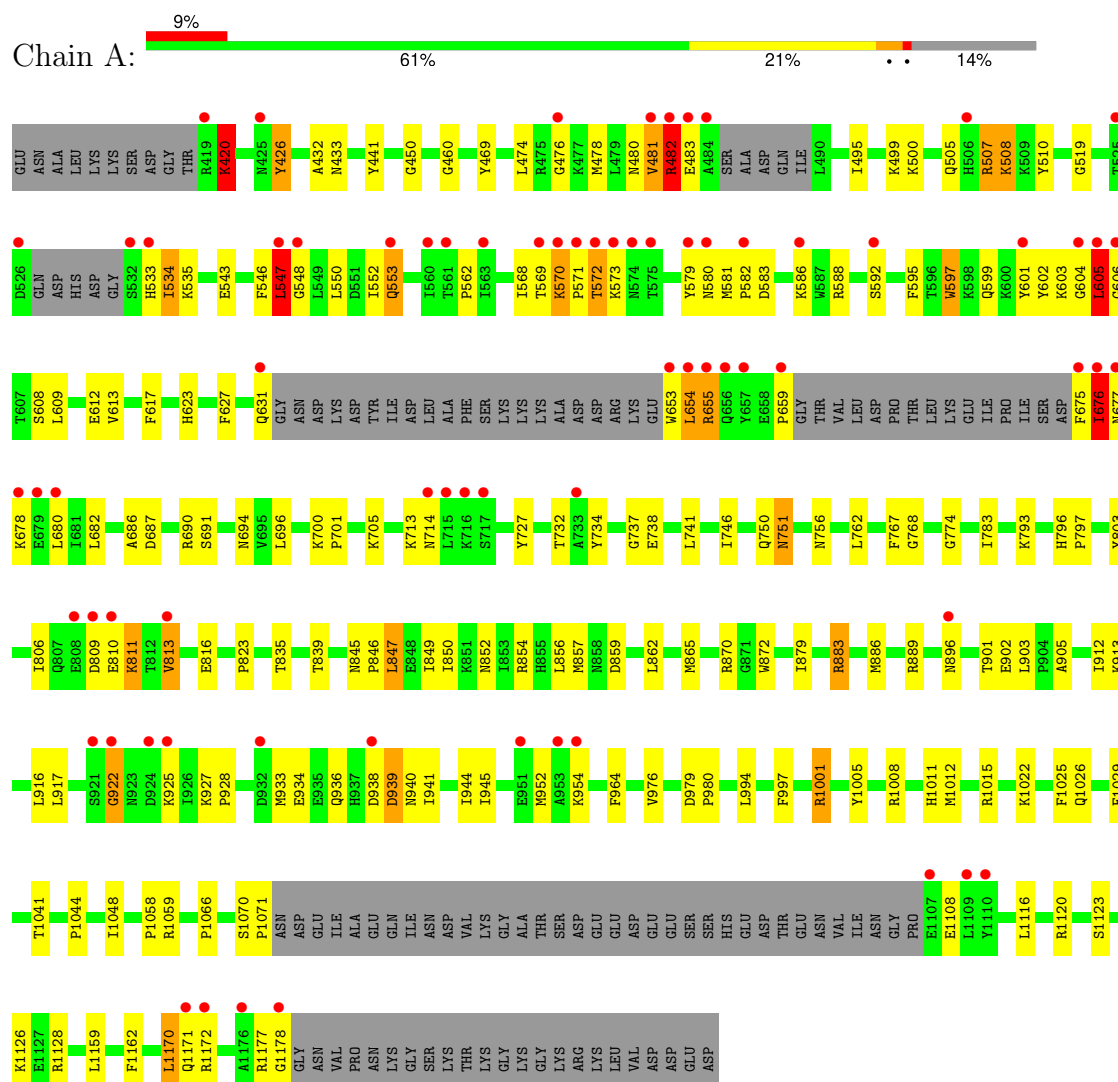
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	217	Total	O	0	0
			217	217		

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: TOPOISOMERASE II



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	103.40Å 145.70Å 161.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.00 – 2.50 14.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.2 (14.00-2.50) 97.3 (14.00-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.58 (at 2.51Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.227 , 0.269 0.236 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	46.7	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5815	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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	0.46	0/5727	0.97	21/7718 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	659	PRO	CA-C-O	10.18	149.55	119.00
1	A	654	LEU	N-CA-C	9.36	123.99	107.80
1	A	426	TYR	CA-C-N	9.01	129.19	119.28
1	A	426	TYR	C-N-CA	9.01	129.19	119.28
1	A	700	LYS	N-CA-C	-8.63	98.06	110.40
1	A	482	ARG	N-CA-C	7.01	119.24	109.15
1	A	816	GLU	N-CA-C	-6.56	104.13	111.28
1	A	691	SER	N-CA-C	5.96	120.73	112.45
1	A	597	TRP	N-CA-C	5.85	118.31	109.41
1	A	793	LYS	N-CA-C	-5.67	106.21	113.23
1	A	676	ILE	N-CA-C	-5.65	97.58	109.34
1	A	569	THR	N-CA-C	5.53	120.91	113.72
1	A	803	TYR	N-CA-C	5.52	117.82	110.53
1	A	927	LYS	CA-C-N	5.47	126.68	119.84
1	A	927	LYS	C-N-CA	5.47	126.68	119.84
1	A	1178	GLY	CA-C-O	5.46	132.27	120.80
1	A	939	ASP	N-CA-C	-5.46	106.62	113.72
1	A	1041	THR	N-CA-C	5.33	118.29	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	938	ASP	N-CA-C	-5.24	99.64	110.80
1	A	605	LEU	N-CA-C	5.23	120.71	110.56
1	A	1108	GLU	N-CA-C	-5.21	107.09	113.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	676	ILE	Mainchain

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	5598	0	5637	158	0
2	A	217	0	0	3	0
All	All	5815	0	5637	158	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 14.

All (158) 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:603:LYS:HB2	1:A:675:PHE:HE1	1.22	1.00
1:A:482:ARG:HD3	1:A:482:ARG:H	1.25	0.98
1:A:627:PHE:HD1	1:A:654:LEU:HD12	1.38	0.88
1:A:623:HIS:CE1	1:A:654:LEU:HD22	2.10	0.87
1:A:903:LEU:HG	1:A:941:ILE:HD13	1.59	0.83
1:A:605:LEU:HG	1:A:606:GLY:N	1.96	0.81
1:A:481:VAL:HG12	1:A:482:ARG:NH1	1.98	0.78
1:A:570:LYS:HB3	1:A:571:PRO:HD3	1.65	0.78
1:A:1120:ARG:HH11	1:A:1120:ARG:HG3	1.49	0.76
1:A:603:LYS:CB	1:A:675:PHE:HE1	1.98	0.75
1:A:568:ILE:O	1:A:573:LYS:HA	1.86	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:603:LYS:HB2	1:A:675:PHE:CE1	2.14	0.73
1:A:655:ARG:HD3	1:A:655:ARG:H	1.53	0.73
1:A:627:PHE:HD1	1:A:654:LEU:CD1	2.02	0.72
1:A:547:LEU:H	1:A:547:LEU:HD22	1.55	0.71
1:A:655:ARG:H	1:A:655:ARG:CD	2.06	0.67
1:A:896:ASN:HA	1:A:952:MET:HE3	1.77	0.67
1:A:481:VAL:HG12	1:A:482:ARG:HH11	1.60	0.66
1:A:852:ASN:HB3	1:A:994:LEU:HD21	1.78	0.65
1:A:588:ARG:HA	1:A:592:SER:OG	1.96	0.65
1:A:849:ILE:CG1	1:A:865:MET:HE1	2.27	0.64
1:A:856:LEU:HG	1:A:862:LEU:HD21	1.79	0.64
1:A:482:ARG:HD3	1:A:482:ARG:N	2.05	0.63
1:A:562:PRO:HA	1:A:579:TYR:HA	1.79	0.63
1:A:997:PHE:CZ	1:A:1001:ARG:HG3	2.34	0.63
1:A:854:ARG:HA	1:A:857:MET:HE2	1.81	0.63
1:A:746:ILE:O	1:A:750:GLN:HG3	1.99	0.63
1:A:768:GLY:HA3	1:A:774:GLY:HA2	1.81	0.62
1:A:849:ILE:HG13	1:A:865:MET:HE1	1.82	0.61
1:A:481:VAL:HG12	1:A:482:ARG:HD3	1.82	0.61
1:A:627:PHE:HA	1:A:654:LEU:HD11	1.82	0.60
1:A:482:ARG:H	1:A:482:ARG:CD	2.01	0.60
1:A:510:TYR:O	1:A:548:GLY:HA3	2.01	0.60
1:A:1044:PRO:O	1:A:1048:ILE:HG13	2.00	0.60
1:A:627:PHE:CD1	1:A:654:LEU:HD12	2.30	0.60
1:A:571:PRO:HG2	1:A:595:PHE:HD1	1.67	0.59
1:A:570:LYS:HE2	1:A:571:PRO:HG3	1.85	0.59
1:A:627:PHE:O	1:A:631:GLN:HG2	2.01	0.59
1:A:568:ILE:HG23	1:A:595:PHE:HB3	1.84	0.58
1:A:507:ARG:HH11	1:A:507:ARG:HG3	1.68	0.58
1:A:823:PRO:HG3	1:A:997:PHE:CD1	2.39	0.58
1:A:474:LEU:HD23	1:A:534:ILE:HD11	1.86	0.58
1:A:553:GLN:N	1:A:553:GLN:HE21	2.01	0.57
1:A:756:ASN:O	1:A:870:ARG:NH1	2.38	0.57
1:A:562:PRO:HG3	1:A:579:TYR:CE2	2.40	0.56
1:A:655:ARG:HD3	1:A:655:ARG:N	2.21	0.56
1:A:922:GLY:H	1:A:928:PRO:HG3	1.71	0.56
1:A:701:PRO:O	1:A:705:LYS:HG3	2.07	0.55
1:A:602:TYR:CZ	1:A:604:GLY:HA3	2.43	0.54
1:A:507:ARG:HG3	1:A:507:ARG:NH1	2.23	0.54
1:A:599:GLN:NE2	1:A:737:GLY:HA2	2.22	0.54
1:A:940:ASN:HD22	1:A:940:ASN:N	2.06	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:623:HIS:HB2	2:A:1363:HOH:O	2.08	0.53
1:A:751:ASN:H	1:A:751:ASN:HD22	1.57	0.53
1:A:687:ASP:OD2	1:A:835:THR:HG23	2.09	0.53
1:A:571:PRO:HG2	1:A:595:PHE:CD1	2.43	0.53
1:A:571:PRO:O	1:A:572:THR:HB	2.09	0.53
1:A:603:LYS:CB	1:A:675:PHE:CE1	2.83	0.53
1:A:627:PHE:CD1	1:A:654:LEU:CD1	2.88	0.52
1:A:903:LEU:CG	1:A:941:ILE:HD13	2.36	0.52
1:A:608:SER:O	1:A:612:GLU:HG2	2.09	0.52
1:A:850:ILE:HD11	1:A:1162:PHE:CD2	2.45	0.51
1:A:806:ILE:O	1:A:813:VAL:HG23	2.09	0.51
1:A:548:GLY:O	1:A:552:ILE:HG13	2.10	0.51
1:A:823:PRO:HG2	1:A:994:LEU:HD12	1.92	0.51
1:A:601:TYR:HE1	1:A:675:PHE:CE1	2.28	0.51
1:A:495:ILE:O	1:A:499:LYS:HG3	2.11	0.51
1:A:1120:ARG:HH11	1:A:1120:ARG:CG	2.23	0.50
1:A:653:TRP:O	1:A:654:LEU:HD23	2.11	0.50
1:A:678:LYS:O	1:A:682:LEU:HB2	2.11	0.49
1:A:592:SER:HB3	1:A:597:TRP:NE1	2.27	0.49
1:A:811:LYS:HB2	1:A:811:LYS:NZ	2.26	0.49
1:A:839:THR:HB	1:A:976:VAL:HB	1.93	0.49
1:A:432:ALA:HB2	1:A:469:TYR:C	2.38	0.49
1:A:605:LEU:HG	1:A:606:GLY:H	1.73	0.49
1:A:1022:LYS:O	1:A:1026:GLN:HG3	2.12	0.49
1:A:823:PRO:HD2	1:A:849:ILE:CG2	2.43	0.48
1:A:583:ASP:O	1:A:586:LYS:HG2	2.14	0.48
1:A:979:ASP:HB2	1:A:980:PRO:CD	2.44	0.48
1:A:879:ILE:HD11	1:A:883:ARG:NH1	2.28	0.48
1:A:925:LYS:HG2	1:A:925:LYS:O	2.14	0.48
1:A:1008:ARG:O	1:A:1012:MET:HG3	2.14	0.48
1:A:845:ASN:C	1:A:845:ASN:HD22	2.22	0.47
1:A:746:ILE:HG23	1:A:767:PHE:CD2	2.49	0.47
1:A:751:ASN:O	1:A:870:ARG:NH2	2.47	0.47
1:A:676:ILE:C	1:A:678:LYS:N	2.73	0.47
1:A:508:LYS:O	1:A:546:PHE:HA	2.13	0.47
1:A:653:TRP:O	1:A:654:LEU:HG	2.14	0.47
1:A:676:ILE:C	1:A:678:LYS:H	2.23	0.47
1:A:1123:SER:HA	1:A:1128:ARG:HG2	1.97	0.47
1:A:734:TYR:CE1	1:A:741:LEU:HD13	2.49	0.47
1:A:823:PRO:HD2	1:A:849:ILE:HG21	1.97	0.46
1:A:852:ASN:HB3	1:A:994:LEU:CD2	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:420:LYS:HE3	1:A:433:ASN:O	2.15	0.46
1:A:850:ILE:HD11	1:A:1162:PHE:CE2	2.51	0.46
1:A:562:PRO:O	1:A:602:TYR:HE1	1.99	0.46
1:A:441:TYR:O	1:A:519:GLY:HA3	2.16	0.46
1:A:694:ASN:ND2	1:A:696:LEU:H	2.13	0.46
1:A:543:GLU:HG3	1:A:550:LEU:HD22	1.97	0.46
1:A:592:SER:HB3	1:A:597:TRP:HE1	1.81	0.45
1:A:872:TRP:CH2	1:A:886:MET:HG3	2.52	0.45
1:A:588:ARG:O	1:A:592:SER:HB2	2.16	0.45
1:A:601:TYR:CE2	1:A:680:LEU:HA	2.51	0.45
1:A:845:ASN:ND2	1:A:847:LEU:H	2.14	0.45
1:A:482:ARG:NH1	1:A:482:ARG:O	2.49	0.45
1:A:1029:PHE:CB	1:A:1116:LEU:HD22	2.47	0.45
1:A:879:ILE:HD11	1:A:883:ARG:HH11	1.82	0.45
1:A:768:GLY:CA	1:A:774:GLY:HA2	2.46	0.44
1:A:1001:ARG:HD3	1:A:1005:TYR:HE2	1.82	0.44
1:A:510:TYR:HD2	1:A:546:PHE:HB3	1.82	0.44
1:A:939:ASP:C	1:A:940:ASN:HD22	2.26	0.44
1:A:478:MET:HG3	1:A:480:ASN:OD1	2.18	0.44
1:A:811:LYS:NZ	1:A:811:LYS:CB	2.80	0.44
1:A:450:GLY:HA2	1:A:476:GLY:O	2.17	0.44
1:A:623:HIS:CE1	1:A:655:ARG:O	2.71	0.44
1:A:767:PHE:CE1	1:A:783:ILE:HB	2.53	0.44
1:A:889:ARG:HB3	1:A:902:GLU:HB3	2.00	0.43
1:A:945:ILE:HD13	1:A:964:PHE:CE1	2.53	0.43
1:A:603:LYS:CG	1:A:605:LEU:HB3	2.49	0.43
1:A:654:LEU:HD23	1:A:655:ARG:CZ	2.49	0.43
1:A:746:ILE:HG23	1:A:767:PHE:HD2	1.83	0.43
1:A:811:LYS:HB2	1:A:811:LYS:HZ2	1.84	0.43
1:A:1025:PHE:CD1	1:A:1058:PRO:HD2	2.54	0.43
1:A:482:ARG:HH11	1:A:482:ARG:N	2.17	0.42
1:A:570:LYS:CE	1:A:571:PRO:HG3	2.48	0.42
1:A:609:LEU:O	1:A:613:VAL:HG23	2.19	0.42
1:A:653:TRP:O	1:A:655:ARG:NH1	2.51	0.42
1:A:845:ASN:ND2	1:A:847:LEU:HB2	2.34	0.42
1:A:796:HIS:HA	1:A:797:PRO:HD2	1.92	0.42
1:A:426:TYR:HE2	1:A:500:LYS:HG2	1.85	0.42
1:A:849:ILE:HG12	1:A:865:MET:HE1	2.00	0.42
1:A:933:MET:SD	1:A:934:GLU:N	2.92	0.42
1:A:581:MET:HB3	1:A:582:PRO:HD3	2.01	0.42
1:A:686:ALA:O	1:A:690:ARG:HG3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:535:LYS:HE3	2:A:1219:HOH:O	2.19	0.42
1:A:653:TRP:O	1:A:654:LEU:CG	2.68	0.42
1:A:460:GLY:HA3	1:A:617:PHE:CD2	2.55	0.42
1:A:510:TYR:HE2	1:A:546:PHE:CD1	2.38	0.42
1:A:912:ILE:O	1:A:916:LEU:HG	2.20	0.42
1:A:936:GLN:NE2	1:A:944:ILE:HD12	2.35	0.42
1:A:676:ILE:O	1:A:678:LYS:N	2.53	0.41
1:A:732:THR:HA	1:A:813:VAL:HG12	2.02	0.41
1:A:713:LYS:HB2	1:A:727:TYR:CE2	2.55	0.41
1:A:1070:SER:HA	1:A:1071:PRO:HD3	1.89	0.41
1:A:886:MET:SD	1:A:905:ALA:HB3	2.61	0.41
1:A:809:ASP:O	1:A:810:GLU:HB2	2.20	0.41
1:A:654:LEU:HD23	1:A:654:LEU:HA	1.90	0.41
1:A:676:ILE:HD13	1:A:676:ILE:HA	2.00	0.41
1:A:676:ILE:HG22	1:A:678:LYS:CG	2.51	0.41
1:A:732:THR:HA	1:A:813:VAL:CG1	2.51	0.41
1:A:845:ASN:HD22	1:A:846:PRO:N	2.19	0.41
1:A:845:ASN:HD21	1:A:847:LEU:HB2	1.86	0.41
1:A:1011:HIS:O	1:A:1015:ARG:HD2	2.21	0.41
1:A:1170:LEU:HD12	1:A:1170:LEU:HA	1.85	0.41
1:A:478:MET:HB2	2:A:1376:HOH:O	2.21	0.40
1:A:751:ASN:HD22	1:A:751:ASN:N	2.19	0.40
1:A:845:ASN:HD22	1:A:847:LEU:H	1.70	0.40
1:A:1059:ARG:O	1:A:1066:PRO:HA	2.22	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	667/793 (84%)	626 (94%)	35 (5%)	6 (1%)	14 27

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	572	THR
1	A	677	ASN
1	A	420	LYS
1	A	922	GLY
1	A	547	LEU
1	A	570	LYS

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	603/701 (86%)	569 (94%)	34 (6%)	19	39

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	420	LYS
1	A	481	VAL
1	A	482	ARG
1	A	483	GLU
1	A	505	GLN
1	A	507	ARG
1	A	508	LYS
1	A	533	HIS
1	A	534	ILE
1	A	547	LEU
1	A	553	GLN
1	A	580	ASN
1	A	605	LEU
1	A	655	ARG
1	A	714	ASN
1	A	738	GLU
1	A	751	ASN
1	A	762	LEU
1	A	811	LYS

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Mol	Chain	Res	Type
1	A	813	VAL
1	A	847	LEU
1	A	859	ASP
1	A	883	ARG
1	A	901	THR
1	A	913	LYS
1	A	917	LEU
1	A	954	LYS
1	A	1001	ARG
1	A	1126	LYS
1	A	1159	LEU
1	A	1170	LEU
1	A	1171	GLN
1	A	1172	ARG
1	A	1177	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	506	HIS
1	A	540	ASN
1	A	553	GLN
1	A	580	ASN
1	A	593	HIS
1	A	623	HIS
1	A	694	ASN
1	A	714	ASN
1	A	735	HIS
1	A	736	HIS
1	A	743	GLN
1	A	751	ASN
1	A	845	ASN
1	A	940	ASN
1	A	974	ASN
1	A	1017	GLN
1	A	1026	GLN
1	A	1130	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](#)

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 [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	679/793 (85%)	0.35	73 (10%) 11 9	23, 50, 99, 100	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	484	ALA	6.2
1	A	676	ILE	5.6
1	A	657	TYR	5.0
1	A	675	PHE	5.0
1	A	1178	GLY	4.9
1	A	809	ASP	4.9
1	A	570	LYS	4.6
1	A	659	PRO	4.4
1	A	653	TRP	4.2
1	A	601	TYR	4.0
1	A	575	THR	3.9
1	A	715	LEU	3.8
1	A	506	HIS	3.7
1	A	553	GLN	3.7
1	A	532	SER	3.7
1	A	606	GLY	3.7
1	A	569	THR	3.6
1	A	678	LYS	3.5
1	A	677	ASN	3.4
1	A	592	SER	3.2
1	A	419	ARG	3.2
1	A	574	ASN	3.2
1	A	655	ARG	3.2
1	A	938	ASP	3.2
1	A	481	VAL	3.1
1	A	571	PRO	3.1
1	A	604	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	476	GLY	3.1
1	A	922	GLY	3.1
1	A	1107	GLU	3.0
1	A	1109	LEU	2.9
1	A	605	LEU	2.9
1	A	547	LEU	2.9
1	A	526	ASP	2.8
1	A	631	GLN	2.8
1	A	679	GLU	2.7
1	A	714	ASN	2.7
1	A	560	ILE	2.6
1	A	810	GLU	2.6
1	A	1176	ALA	2.6
1	A	533	HIS	2.6
1	A	1172	ARG	2.5
1	A	654	LEU	2.5
1	A	953	ALA	2.5
1	A	525	THR	2.5
1	A	572	THR	2.5
1	A	717	SER	2.5
1	A	548	GLY	2.5
1	A	716	LYS	2.5
1	A	925	LYS	2.5
1	A	425	ASN	2.5
1	A	579	TYR	2.4
1	A	1110	TYR	2.4
1	A	656	GLN	2.4
1	A	1171	GLN	2.3
1	A	921	SER	2.3
1	A	680	LEU	2.3
1	A	932	ASP	2.3
1	A	733	ALA	2.2
1	A	896	ASN	2.2
1	A	582	PRO	2.2
1	A	483	GLU	2.2
1	A	482	ARG	2.2
1	A	573	LYS	2.2
1	A	586	LYS	2.2
1	A	561	THR	2.1
1	A	808	GLU	2.1
1	A	951	GLU	2.1
1	A	580	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	563	ILE	2.0
1	A	813	VAL	2.0
1	A	954	LYS	2.0
1	A	924	ASP	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.