



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:45 AM UTC

PDB ID : 3K4X / pdb_00003k4x
Title : Eukaryotic Sliding Clamp PCNA Bound to DNA
Authors : McNally, R.; Kuriyan, J.
Deposited on : 2009-10-06
Resolution : 2.98 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

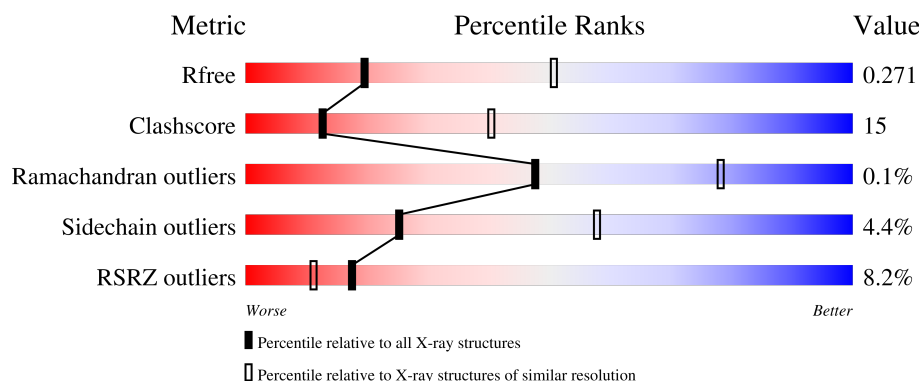
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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	798	7% 67% 26% • 5%
2	B	10	40% 100%
3	C	14	29% 71% 29%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6354 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 Proliferating cell nuclear antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	758	5947	3802	939	1179	27	0	0	0

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP P15873
A	-4	PRO	-	expression tag	UNP P15873
A	-3	HIS	-	expression tag	UNP P15873
A	-2	MET	-	expression tag	UNP P15873
A	-1	ALA	-	expression tag	UNP P15873
A	0	SER	-	expression tag	UNP P15873
A	258	GLY	-	linker	UNP P15873
A	259	SER	-	linker	UNP P15873
A	260	ASN	-	linker	UNP P15873
A	261	SER	-	linker	UNP P15873
A	262	GLN	-	linker	UNP P15873
A	263	SER	-	linker	UNP P15873
A	264	ASN	-	linker	UNP P15873
A	265	GLY	-	linker	UNP P15873
A	266	SER	-	linker	UNP P15873
A	267	GLY	-	linker	UNP P15873
A	268	ALA	-	linker	UNP P15873
A	258	GLY	-	linker	UNP P15873
A	259	SER	-	linker	UNP P15873
A	260	ASN	-	linker	UNP P15873
A	261	SER	-	linker	UNP P15873
A	262	GLN	-	linker	UNP P15873
A	263	SER	-	linker	UNP P15873
A	264	ASN	-	linker	UNP P15873
A	265	GLY	-	linker	UNP P15873
A	266	SER	-	linker	UNP P15873
A	267	GLY	-	linker	UNP P15873

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Chain	Residue	Modelled	Actual	Comment	Reference
A	268	ALA	-	linker	UNP P15873

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*CP*CP*AP*TP*CP*GP*TP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	10	Total	C	N	O	P	0	0	0
			197	96	33	59	9			

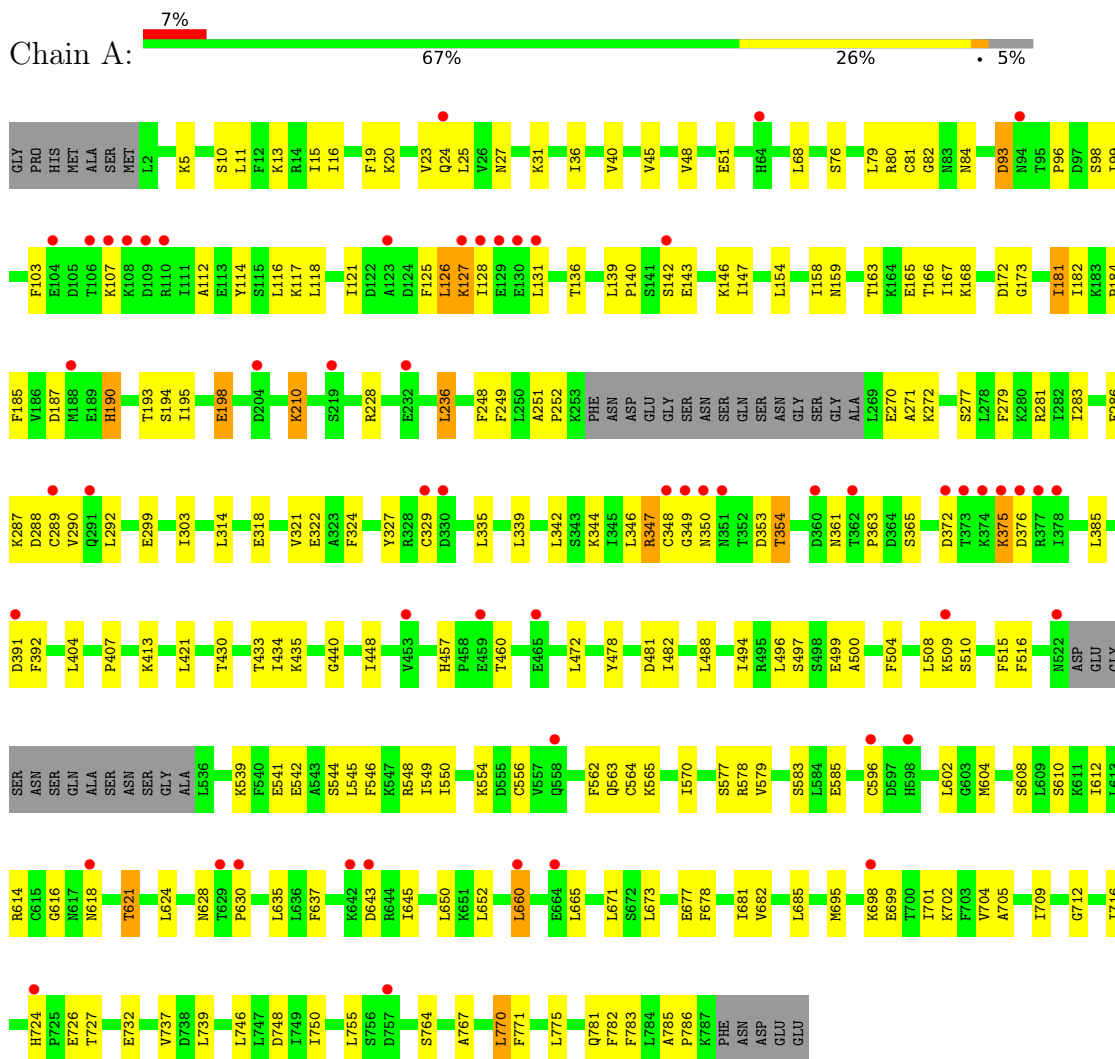
- Molecule 3 is a DNA chain called DNA (5'-D(*TP*TP*TP*TP*AP*TP*AP*CP*GP*AP*TP*GP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	P	0	0	0
			210	99	42	59	10			

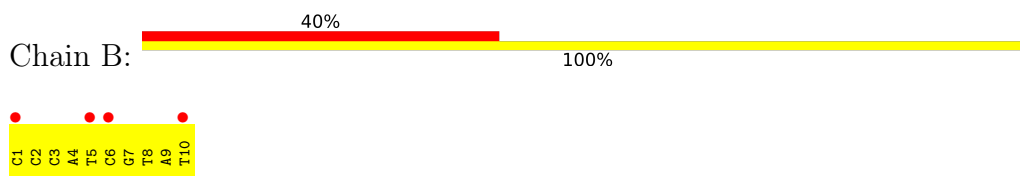
3 Residue-property plots

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: Proliferating cell nuclear antigen



• Molecule 2: DNA (5'-D(*CP*CP*CP*AP*TP*CP*GP*TP*AP*T)-3')



- Molecule 3: DNA (5'-D(*TP*TP*TP*TP*AP*TP*AP*CP*GP*AP*TP*GP*GP*G)-3')

Chain C: 

DT	DT	DT	DT	A5	T6	A7	C8	G9	A10	T11	G12	G13	G14
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.94Å 93.71Å 136.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.86 – 2.98 46.86 – 2.99	Depositor EDS
% Data completeness (in resolution range)	97.6 (46.86-2.98) 97.6 (46.86-2.99)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.4_162	Depositor
R, R_{free}	0.223 , 0.280 0.214 , 0.271	Depositor DCC
R_{free} test set	1121 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	52.4	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 52.9	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	0.91	EDS
Total number of atoms	6354	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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	0.25	0/6035	0.69	0/8142
2	B	0.45	0/219	1.02	0/335
3	C	0.47	0/236	0.96	0/363
All	All	0.27	0/6490	0.72	0/8840

There are no bond length outliers.

There are no bond angle outliers.

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	5947	0	6009	155	0
2	B	197	0	115	20	0
3	C	210	0	113	18	0
All	All	6354	0	6237	191	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 15.

All (191) 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:142:SER:HA	2:B:6:DC:OP1	1.85	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:544:SER:HB3	1:A:618:ASN:HB3	1.69	0.74
1:A:24:GLN:HG3	1:A:25:LEU:HG	1.71	0.73
1:A:347:ARG:HA	1:A:347:ARG:HE	1.54	0.72
1:A:190:HIS:O	1:A:193:THR:HG22	1.90	0.72
1:A:539:LYS:HE3	1:A:621:THR:HG21	1.74	0.70
1:A:127:LYS:C	1:A:128:ILE:HD12	2.17	0.70
1:A:10:SER:HB3	1:A:84:ASN:HB3	1.74	0.69
1:A:303:ILE:HG22	1:A:318:GLU:HG2	1.75	0.68
1:A:457:HIS:O	1:A:460:THR:HG22	1.94	0.68
1:A:272:LYS:HE3	1:A:354:THR:HG21	1.77	0.66
1:A:671:LEU:HD13	1:A:701:ILE:HD13	1.75	0.66
1:A:270:GLU:O	1:A:270:GLU:HG3	1.97	0.65
1:A:755:LEU:HB3	1:A:775:LEU:HD21	1.79	0.65
1:A:146:LYS:HZ1	1:A:349:GLY:HA2	1.62	0.64
1:A:361:ASN:O	1:A:363:PRO:HD3	1.97	0.64
1:A:544:SER:HB2	1:A:548:ARG:HH12	1.63	0.64
1:A:563:GLN:HB3	1:A:565:LYS:NZ	2.13	0.63
1:A:182:ILE:HD12	1:A:195:ILE:HD13	1.81	0.62
1:A:724:HIS:O	1:A:727:THR:HG22	1.99	0.62
1:A:314:LEU:HB3	1:A:516:PHE:HB2	1.82	0.61
1:A:365:SER:HA	1:A:385:LEU:HG	1.83	0.60
1:A:404:LEU:HD13	1:A:434:ILE:HD13	1.82	0.60
1:A:277:SER:HB2	1:A:281:ARG:NH1	2.18	0.59
1:A:76:SER:O	1:A:80:ARG:HG3	2.02	0.59
1:A:131:LEU:H	1:A:131:LEU:HD23	1.68	0.59
1:A:562:PHE:CD2	1:A:604:MET:HE2	2.39	0.58
1:A:286:PHE:HB3	1:A:339:LEU:HD13	1.87	0.57
1:A:287:LYS:O	1:A:288:ASP:HB2	2.04	0.57
1:A:660:LEU:H	1:A:660:LEU:HD22	1.68	0.57
3:C:8:DC:H2''	3:C:9:DG:OP2	2.05	0.57
2:B:5:DT:H2''	2:B:6:DC:OP2	2.05	0.57
2:B:8:DT:H2''	2:B:9:DA:OP2	2.05	0.56
3:C:5:DA:H2''	3:C:6:DT:OP2	2.05	0.56
2:B:7:DG:H1'	2:B:8:DT:H5'	1.88	0.56
3:C:10:DA:H2''	3:C:11:DT:OP2	2.05	0.56
3:C:10:DA:H1'	3:C:11:DT:H5'	1.88	0.56
3:C:13:DG:H1'	3:C:14:DG:H5'	1.88	0.56
2:B:4:DA:H1'	2:B:5:DT:H5'	1.88	0.56
3:C:6:DT:H1'	3:C:7:DA:H5'	1.88	0.56
2:B:1:DC:H2''	2:B:2:DC:OP2	2.05	0.56
2:B:2:DC:H1'	2:B:3:DC:H5'	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:7:DA:H2''	3:C:8:DC:OP2	2.05	0.56
3:C:9:DG:H1'	3:C:10:DA:H5'	1.87	0.56
3:C:9:DG:H2''	3:C:10:DA:OP2	2.05	0.56
2:B:4:DA:H2''	2:B:5:DT:OP2	2.05	0.56
3:C:12:DG:H2''	3:C:13:DG:OP2	2.05	0.56
3:C:13:DG:H2''	3:C:14:DG:OP2	2.05	0.56
3:C:8:DC:H1'	3:C:9:DG:H5'	1.88	0.56
2:B:1:DC:H1'	2:B:2:DC:H5'	1.88	0.56
3:C:7:DA:H1'	3:C:8:DC:H5'	1.88	0.56
1:A:136:THR:HB	1:A:198:GLU:HG2	1.88	0.56
2:B:3:DC:H1'	2:B:4:DA:H5'	1.88	0.56
3:C:5:DA:H1'	3:C:6:DT:H5'	1.88	0.56
3:C:11:DT:H1'	3:C:12:DG:H5'	1.87	0.56
1:A:68:LEU:HD23	1:A:99:ILE:HD12	1.87	0.56
2:B:5:DT:H1'	2:B:6:DC:H5'	1.88	0.56
2:B:6:DC:H2''	2:B:7:DG:OP2	2.05	0.56
1:A:570:ILE:HG22	1:A:585:GLU:HG2	1.88	0.55
1:A:610:SER:O	1:A:614:ARG:HG3	2.06	0.55
2:B:9:DA:H1'	2:B:10:DT:H5'	1.88	0.55
3:C:6:DT:H2''	3:C:7:DA:OP2	2.05	0.55
1:A:93:ASP:OD2	1:A:93:ASP:N	2.39	0.55
1:A:347:ARG:HE	1:A:347:ARG:CA	2.18	0.55
2:B:3:DC:H2''	2:B:4:DA:OP2	2.05	0.55
2:B:7:DG:H2''	2:B:8:DT:OP2	2.05	0.55
2:B:2:DC:H2''	2:B:3:DC:OP2	2.05	0.55
2:B:6:DC:H1'	2:B:7:DG:H5'	1.88	0.55
2:B:9:DA:H2''	2:B:10:DT:OP2	2.05	0.55
3:C:11:DT:H2''	3:C:12:DG:OP2	2.05	0.55
1:A:125:PHE:HE2	1:A:127:LYS:HD2	1.72	0.55
1:A:556:CYS:SG	1:A:748:ASP:HB3	2.46	0.55
2:B:8:DT:H1'	2:B:9:DA:H5'	1.88	0.55
1:A:279:PHE:HD2	1:A:346:LEU:HD11	1.72	0.55
1:A:541:GLU:HG3	1:A:542:GLU:HG2	1.88	0.55
1:A:407:PRO:HG3	1:A:460:THR:HA	1.89	0.54
1:A:286:PHE:HD2	1:A:290:VAL:HG21	1.72	0.54
3:C:12:DG:H1'	3:C:13:DG:H5'	1.88	0.54
1:A:168:LYS:HG3	1:A:181:ILE:HD13	1.88	0.54
1:A:602:LEU:HD12	1:A:652:LEU:HD11	1.89	0.54
1:A:127:LYS:N	1:A:128:ILE:HD12	2.24	0.53
1:A:549:ILE:HA	1:A:755:LEU:HD21	1.90	0.53
1:A:544:SER:O	1:A:548:ARG:HG2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:612:ILE:HG21	1:A:635:LEU:HD13	1.91	0.53
1:A:272:LYS:HE3	1:A:354:THR:CG2	2.39	0.52
1:A:746:LEU:O	1:A:750:ILE:HG23	2.09	0.52
1:A:353:ASP:OD2	1:A:372:ASP:HA	2.10	0.52
1:A:321:VAL:HG13	1:A:327:TYR:HB3	1.91	0.51
1:A:342:LEU:O	1:A:346:LEU:HD23	2.10	0.51
1:A:168:LYS:HG3	1:A:181:ILE:CD1	2.39	0.51
1:A:27:ASN:OD1	1:A:121:ILE:HB	2.10	0.51
1:A:210:LYS:O	1:A:210:LYS:HD3	2.11	0.51
1:A:550:ILE:O	1:A:554:LYS:HB3	2.10	0.51
1:A:299:GLU:H	1:A:299:GLU:CD	2.19	0.51
1:A:322:GLU:HB2	1:A:510:SER:HB2	1.92	0.51
1:A:544:SER:HB2	1:A:548:ARG:NH1	2.26	0.50
1:A:539:LYS:HE3	1:A:621:THR:CG2	2.41	0.50
1:A:126:LEU:C	1:A:128:ILE:HD12	2.37	0.50
1:A:628:ASN:O	1:A:630:PRO:HD3	2.11	0.50
1:A:154:LEU:O	1:A:173:GLY:HA3	2.12	0.49
1:A:116:LEU:HD12	1:A:709:ILE:O	2.13	0.49
1:A:494:ILE:HG22	1:A:496:LEU:HG	1.94	0.49
1:A:127:LYS:HA	1:A:127:LYS:HE3	1.94	0.49
1:A:478:TYR:O	1:A:482:ILE:HG13	2.13	0.49
1:A:677:GLU:O	1:A:681:ILE:HG13	2.13	0.49
1:A:185:PHE:HE1	1:A:187:ASP:HB2	1.78	0.48
1:A:771:PHE:HB2	1:A:782:PHE:HB2	1.95	0.48
1:A:114:TYR:HA	1:A:712:GLY:HA2	1.95	0.48
1:A:112:ALA:HB1	1:A:114:TYR:CE1	2.48	0.48
1:A:136:THR:HG23	1:A:228:ARG:HG2	1.96	0.48
1:A:322:GLU:HB2	1:A:510:SER:CB	2.44	0.48
1:A:314:LEU:HD23	1:A:516:PHE:CD1	2.49	0.47
1:A:13:LYS:HE2	1:A:84:ASN:HD21	1.79	0.47
1:A:93:ASP:O	1:A:96:PRO:HG3	2.14	0.47
1:A:430:THR:OG1	1:A:433:THR:HB	2.15	0.47
1:A:19:PHE:O	1:A:23:VAL:HG22	2.13	0.47
1:A:128:ILE:HD12	1:A:128:ILE:N	2.30	0.47
1:A:702:LYS:HE2	1:A:704:VAL:HG21	1.96	0.47
1:A:286:PHE:O	1:A:290:VAL:HG22	2.14	0.47
1:A:724:HIS:CE1	1:A:726:GLU:HB2	2.49	0.47
1:A:142:SER:CA	2:B:6:DC:OP1	2.60	0.47
1:A:350:ASN:HB2	1:A:353:ASP:CG	2.40	0.47
1:A:546:PHE:O	1:A:550:ILE:HG12	2.15	0.47
1:A:563:GLN:NE2	1:A:570:ILE:HD11	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:SER:HB2	1:A:116:LEU:O	2.14	0.46
1:A:103:PHE:HB2	1:A:112:ALA:HB3	1.96	0.46
1:A:112:ALA:HB1	1:A:114:TYR:HE1	1.80	0.46
1:A:81:CYS:HB3	1:A:114:TYR:OH	2.16	0.46
1:A:375:LYS:HE3	1:A:375:LYS:HA	1.98	0.46
1:A:143:GLU:O	1:A:147:ILE:HG13	2.15	0.46
1:A:472:LEU:HD11	1:A:497:SER:O	2.15	0.46
1:A:770:LEU:HG	1:A:783:PHE:CE2	2.51	0.46
1:A:344:LYS:O	1:A:347:ARG:HB2	2.16	0.45
1:A:139:LEU:H	1:A:139:LEU:HD23	1.81	0.45
1:A:660:LEU:HD22	1:A:660:LEU:N	2.31	0.45
1:A:695:MET:HE3	1:A:702:LYS:HD3	1.97	0.45
1:A:167:ILE:HB	1:A:182:ILE:CG1	2.47	0.45
1:A:187:ASP:OD2	1:A:190:HIS:HD2	2.00	0.45
1:A:497:SER:HB3	1:A:500:ALA:HB3	1.98	0.45
1:A:5:LYS:O	1:A:5:LYS:HG3	2.15	0.45
1:A:277:SER:HB2	1:A:281:ARG:HH11	1.81	0.45
1:A:96:PRO:HB2	1:A:118:LEU:HD12	1.99	0.45
1:A:685:LEU:HB3	1:A:705:ALA:HB2	1.98	0.45
1:A:737:VAL:HG12	1:A:739:LEU:HG	1.99	0.44
1:A:545:LEU:O	1:A:549:ILE:HG13	2.17	0.44
1:A:48:VAL:HG22	1:A:248:PHE:CD2	2.53	0.44
1:A:643:ASP:O	1:A:645:ILE:HG13	2.17	0.44
1:A:583:SER:HB3	1:A:781:GLN:HB2	1.98	0.44
1:A:125:PHE:CE2	1:A:127:LYS:HD2	2.53	0.44
1:A:163:THR:OG1	1:A:166:THR:HB	2.17	0.44
1:A:698:LYS:HG2	1:A:699:GLU:HG3	2.00	0.44
1:A:125:PHE:HE2	1:A:127:LYS:HB2	1.82	0.44
1:A:435:LYS:HG3	1:A:448:ILE:HG12	2.00	0.44
1:A:251:ALA:HA	1:A:252:PRO:HD3	1.88	0.44
1:A:146:LYS:NZ	1:A:349:GLY:HA2	2.31	0.43
1:A:509:LYS:O	1:A:510:SER:HB3	2.17	0.43
1:A:702:LYS:HE2	1:A:704:VAL:CG2	2.48	0.43
1:A:16:ILE:O	1:A:20:LYS:HB3	2.17	0.43
1:A:785:ALA:HA	1:A:786:PRO:HD3	1.87	0.43
1:A:391:ASP:HA	1:A:392:PHE:HA	1.64	0.43
1:A:289:CYS:HB3	1:A:481:ASP:HB3	2.00	0.43
1:A:624:LEU:CD2	1:A:635:LEU:HD23	2.49	0.43
1:A:36:ILE:HG22	1:A:51:GLU:HG2	2.01	0.43
1:A:125:PHE:CE2	1:A:127:LYS:HB2	2.53	0.42
1:A:488:LEU:HD13	1:A:508:LEU:HD21	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:LYS:HE2	1:A:107:LYS:HB3	1.86	0.42
1:A:504:PHE:HB2	1:A:515:PHE:HB2	2.00	0.42
1:A:11:LEU:O	1:A:15:ILE:HG13	2.19	0.42
1:A:140:PRO:HG3	1:A:193:THR:HA	2.02	0.42
1:A:271:ALA:HB1	1:A:324:PHE:CE1	2.54	0.42
1:A:347:ARG:HA	1:A:347:ARG:NE	2.30	0.42
1:A:81:CYS:HB3	1:A:114:TYR:CZ	2.55	0.42
1:A:764:SER:HB3	1:A:767:ALA:HB3	2.00	0.42
1:A:165:GLU:HA	1:A:184:PRO:HG2	2.02	0.41
1:A:350:ASN:HB2	1:A:353:ASP:OD2	2.21	0.41
1:A:187:ASP:H	1:A:194:SER:HB3	1.85	0.41
1:A:286:PHE:HB3	1:A:339:LEU:CD1	2.51	0.41
1:A:421:LEU:O	1:A:440:GLY:HA3	2.19	0.41
1:A:678:PHE:O	1:A:682:VAL:HG23	2.20	0.41
1:A:25:LEU:HB3	1:A:121:ILE:HD13	2.02	0.41
1:A:31:LYS:HB2	1:A:31:LYS:HE3	1.85	0.41
1:A:616:GLY:HA3	1:A:637:PHE:CD2	2.55	0.41
1:A:577:SER:O	1:A:578:ARG:HB2	2.20	0.41
1:A:158:ILE:HG22	1:A:159:ASN:N	2.36	0.40
1:A:279:PHE:O	1:A:283:ILE:HG12	2.21	0.40
1:A:701:ILE:HB	1:A:716:ILE:CG1	2.51	0.40
1:A:236:LEU:HG	1:A:249:PHE:CE2	2.56	0.40
1:A:608:SER:HB3	1:A:650:LEU:HD11	2.03	0.40
1:A:673:LEU:HD23	1:A:673:LEU:N	2.36	0.40
1:A:140:PRO:C	1:A:142:SER:N	2.79	0.40
1:A:564:CYS:O	1:A:565:LYS:HD3	2.20	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	752/798 (94%)	730 (97%)	21 (3%)	1 (0%)	48 78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	82	GLY

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	683/713 (96%)	653 (96%)	30 (4%)	25 58

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	VAL
1	A	45	VAL
1	A	79	LEU
1	A	93	ASP
1	A	117	LYS
1	A	126	LEU
1	A	127	LYS
1	A	172	ASP
1	A	181	ILE
1	A	190	HIS
1	A	198	GLU
1	A	210	LYS
1	A	236	LEU
1	A	292	LEU
1	A	329	CYS
1	A	335	LEU
1	A	347	ARG
1	A	348	CYS
1	A	354	THR
1	A	375	LYS

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Mol	Chain	Res	Type
1	A	376	ASP
1	A	413	LYS
1	A	499	GLU
1	A	579	VAL
1	A	596	CYS
1	A	621	THR
1	A	660	LEU
1	A	665	LEU
1	A	732	GLU
1	A	770	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	29	GLN
1	A	132	GLN
1	A	153	GLN
1	A	190	HIS
1	A	563	GLN
1	A	781	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 [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	758/798 (94%)	0.31	56 (7%)	20 13	15, 46, 111, 210	0
2	B	10/10 (100%)	2.20	4 (40%)	1 1	179, 293, 361, 367	0
3	C	10/14 (71%)	1.98	4 (40%)	1 1	223, 275, 323, 336	0
All	All	778/822 (94%)	0.36	64 (8%)	17 11	15, 47, 131, 367	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	660	LEU	5.9
1	A	127	LYS	5.8
1	A	348	CYS	5.5
1	A	349	GLY	5.2
1	A	291	GLN	4.9
1	A	351	ASN	4.9
1	A	391	ASP	4.8
1	A	204	ASP	4.7
1	A	109	ASP	4.5
1	A	664	GLU	4.5
2	B	1	DC	4.4
1	A	131	LEU	4.3
1	A	110	ARG	4.0
1	A	377	ARG	3.7
1	A	376	ASP	3.7
1	A	104	GLU	3.6
1	A	329	CYS	3.6
1	A	128	ILE	3.4
1	A	378	ILE	3.4
1	A	94	ASN	3.3
3	C	14	DG	3.1
1	A	375	LYS	3.0
1	A	643	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	6	DC	3.0
1	A	64	HIS	3.0
1	A	289	CYS	2.9
1	A	142	SER	2.9
1	A	129	GLU	2.9
1	A	372	ASP	2.8
1	A	465	GLU	2.7
1	A	350	ASN	2.6
1	A	360	ASP	2.6
1	A	188	MET	2.6
1	A	373	THR	2.6
1	A	108	LYS	2.5
2	B	10	DT	2.5
1	A	596	CYS	2.5
1	A	629	THR	2.5
1	A	123	ALA	2.5
1	A	642	LYS	2.5
1	A	598	HIS	2.4
1	A	724	HIS	2.4
1	A	630	PRO	2.4
1	A	558	GLN	2.4
1	A	522	ASN	2.3
3	C	5	DA	2.3
3	C	11	DT	2.3
1	A	453	VAL	2.3
1	A	24	GLN	2.2
1	A	698	LYS	2.2
1	A	106	THR	2.2
1	A	107	LYS	2.2
1	A	618	ASN	2.2
1	A	374	LYS	2.2
1	A	757	ASP	2.2
2	B	5	DT	2.1
1	A	219	SER	2.1
3	C	8	DC	2.1
1	A	459	GLU	2.1
1	A	330	ASP	2.1
1	A	362	THR	2.0
1	A	130	GLU	2.0
1	A	232	GLU	2.0
1	A	509	LYS	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.