



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:10 PM UTC

PDB ID : 9DHU / pdb_00009dhu
Title : The Retinoblastoma Protein with Mutation E533K
Authors : Ruiz-Rivera, A.; Castro, A.; Burke, J.R.
Deposited on : 2024-09-04
Resolution : 2.16 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

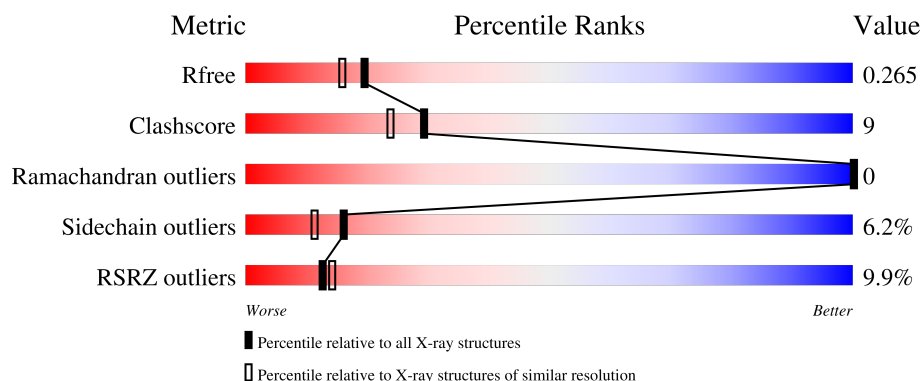
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	390	
1	B	390	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11521 atoms, of which 5768 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 Retinoblastoma-associated protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	336	Total	C	H	N	O	S	0	0	0
			5605	1793	2831	466	495	20			
1	B	353	Total	C	H	N	O	S	0	0	0
			5842	1872	2937	488	524	21			

There are 82 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	377	GLY	-	expression tag	UNP P06400
A	378	GLU	-	expression tag	UNP P06400
A	379	PHE	-	expression tag	UNP P06400
A	533	LYS	GLU	conflict	UNP P06400
A	608	GLU	SER	conflict	UNP P06400
A	612	ALA	SER	conflict	UNP P06400
A	?	-	LYS	deletion	UNP P06400
A	?	-	GLY	deletion	UNP P06400
A	?	-	SER	deletion	UNP P06400
A	?	-	THR	deletion	UNP P06400
A	?	-	THR	deletion	UNP P06400
A	?	-	ARG	deletion	UNP P06400
A	?	-	VAL	deletion	UNP P06400
A	?	-	ASN	deletion	UNP P06400
A	?	-	SER	deletion	UNP P06400
A	?	-	THR	deletion	UNP P06400
A	?	-	ALA	deletion	UNP P06400
A	?	-	ASN	deletion	UNP P06400
A	?	-	ALA	deletion	UNP P06400
A	?	-	GLU	deletion	UNP P06400
A	?	-	THR	deletion	UNP P06400
A	?	-	GLN	deletion	UNP P06400
A	?	-	ALA	deletion	UNP P06400
A	?	-	THR	deletion	UNP P06400
A	?	-	SER	deletion	UNP P06400

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ALA	deletion	UNP P06400
A	?	-	PHE	deletion	UNP P06400
A	?	-	GLN	deletion	UNP P06400
A	?	-	THR	deletion	UNP P06400
A	?	-	GLN	deletion	UNP P06400
A	?	-	LYS	deletion	UNP P06400
A	?	-	PRO	deletion	UNP P06400
A	?	-	LEU	deletion	UNP P06400
A	780	ALA	SER	conflict	UNP P06400
A	787	PHE	-	expression tag	UNP P06400
A	788	MET	-	expression tag	UNP P06400
A	789	GLN	-	expression tag	UNP P06400
A	790	ARG	-	expression tag	UNP P06400
A	791	LEU	-	expression tag	UNP P06400
A	792	LYS	-	expression tag	UNP P06400
A	793	THR	-	expression tag	UNP P06400
B	377	GLY	-	expression tag	UNP P06400
B	378	GLU	-	expression tag	UNP P06400
B	379	PHE	-	expression tag	UNP P06400
B	533	LYS	GLU	conflict	UNP P06400
B	608	GLU	SER	conflict	UNP P06400
B	639	ALA	SER	conflict	UNP P06400
B	?	-	LYS	deletion	UNP P06400
B	?	-	GLY	deletion	UNP P06400
B	?	-	SER	deletion	UNP P06400
B	?	-	THR	deletion	UNP P06400
B	?	-	THR	deletion	UNP P06400
B	?	-	ARG	deletion	UNP P06400
B	?	-	VAL	deletion	UNP P06400
B	?	-	ASN	deletion	UNP P06400
B	?	-	SER	deletion	UNP P06400
B	?	-	THR	deletion	UNP P06400
B	?	-	ALA	deletion	UNP P06400
B	?	-	ASN	deletion	UNP P06400
B	?	-	ALA	deletion	UNP P06400
B	?	-	GLU	deletion	UNP P06400
B	?	-	THR	deletion	UNP P06400
B	?	-	GLN	deletion	UNP P06400
B	?	-	ALA	deletion	UNP P06400
B	?	-	THR	deletion	UNP P06400
B	?	-	SER	deletion	UNP P06400
B	?	-	ALA	deletion	UNP P06400

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	PHE	deletion	UNP P06400
B	?	-	GLN	deletion	UNP P06400
B	?	-	THR	deletion	UNP P06400
B	?	-	GLN	deletion	UNP P06400
B	?	-	LYS	deletion	UNP P06400
B	?	-	PRO	deletion	UNP P06400
B	?	-	LEU	deletion	UNP P06400
B	780	ALA	SER	conflict	UNP P06400
B	787	PHE	-	expression tag	UNP P06400
B	788	MET	-	expression tag	UNP P06400
B	789	GLN	-	expression tag	UNP P06400
B	790	ARG	-	expression tag	UNP P06400
B	791	LEU	-	expression tag	UNP P06400
B	792	LYS	-	expression tag	UNP P06400
B	793	THR	-	expression tag	UNP P06400

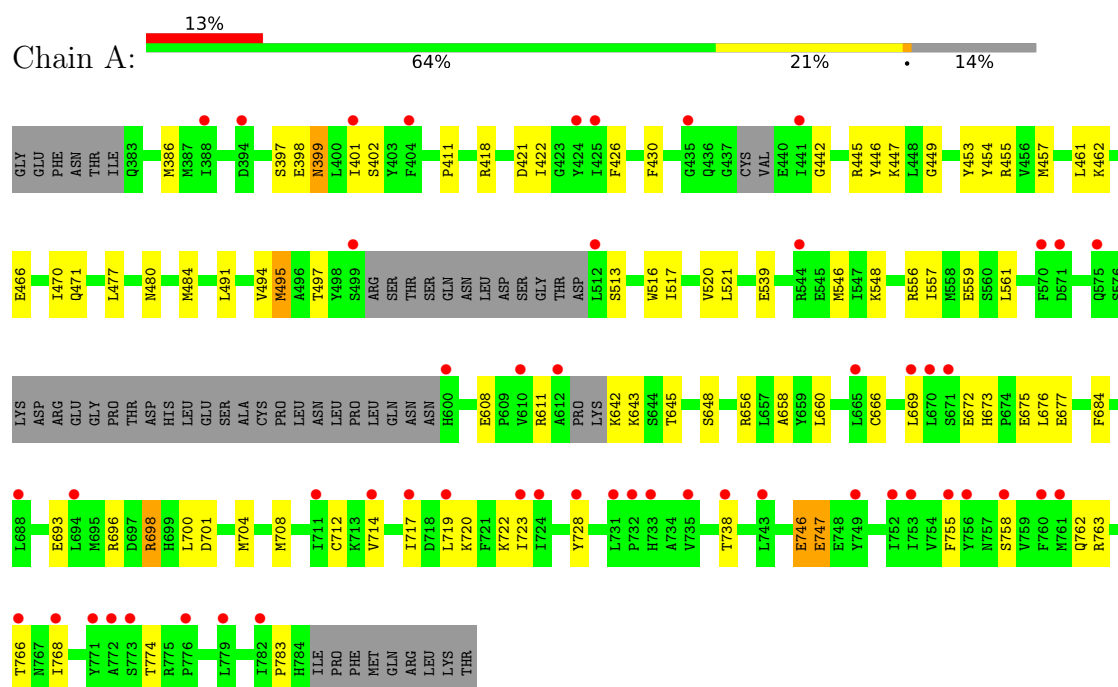
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	9	Total O 9 9	0	0
2	B	65	Total O 65 65	0	0

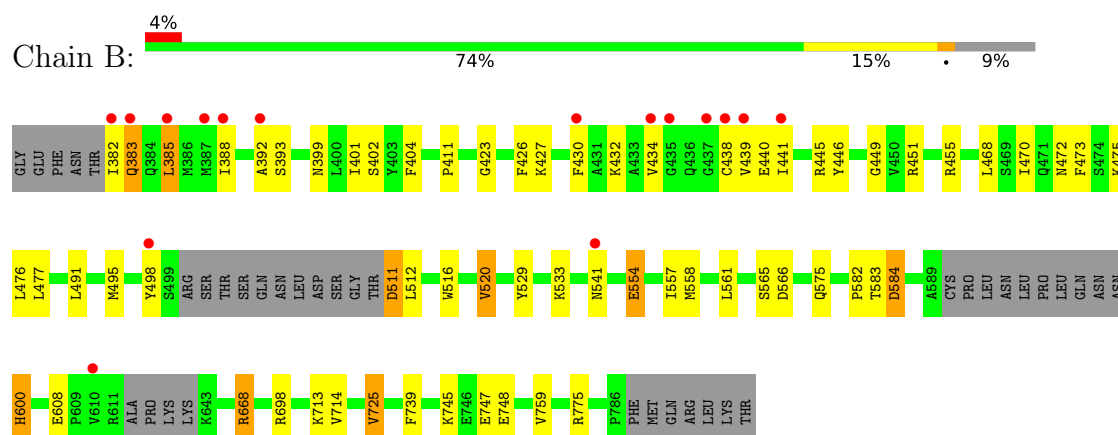
3 Residue-property plots

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: Retinoblastoma-associated protein



• Molecule 1: Retinoblastoma-associated protein



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	250.14Å 250.14Å 35.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.27 – 2.16 47.27 – 2.16	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.27-2.16) 99.9 (47.27-2.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.16Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660, PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.226 , 0.263 0.227 , 0.265	Depositor DCC
R_{free} test set	1996 reflections (4.52%)	wwPDB-VP
Wilson B-factor (Å ²)	38.6	Xtriage
Anisotropy	0.291	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.010 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11521	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% 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	0.49	0/2829	0.75	1/3806 (0.0%)
1	B	0.94	1/2964 (0.0%)	1.01	2/3993 (0.1%)
All	All	0.75	1/5793 (0.0%)	0.90	3/7799 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	582	PRO	C-O	-5.46	1.17	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	696	ARG	CA-C-O	-7.41	113.92	120.88
1	B	383	GLN	N-CA-C	6.45	121.07	112.30
1	B	554	GLU	N-CA-C	-5.15	105.56	111.07

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	2774	2831	2834	60	0
1	B	2905	2937	2955	39	0
2	A	9	0	0	0	0
2	B	65	0	0	2	0
All	All	5753	5768	5789	98	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 9.

All (98) 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:554:GLU:HG2	1:B:558:MET:HE2	1.76	0.68
1:A:666:CYS:SG	1:A:708:MET:HE3	2.36	0.65
1:A:669:LEU:HD22	1:A:723:ILE:HG22	1.79	0.65
1:B:725:VAL:HG13	1:B:739:PHE:CZ	2.34	0.62
1:B:430:PHE:CE1	1:B:434:VAL:HG11	2.35	0.61
1:B:382:ILE:O	1:B:382:ILE:HG22	2.01	0.60
1:A:386:MET:HE2	1:A:497:THR:HG21	1.86	0.58
1:A:701:ASP:HA	1:A:704:MET:HE2	1.87	0.57
1:B:430:PHE:HE1	1:B:434:VAL:HG11	1.68	0.57
1:B:438:CYS:HB3	1:B:441:ILE:HD12	1.85	0.57
1:B:385:LEU:HB2	1:B:541:ASN:HB3	1.88	0.56
1:A:746:GLU:HA	1:A:746:GLU:OE1	2.05	0.56
1:B:600:HIS:CE1	2:B:804:HOH:O	2.58	0.56
1:A:461:LEU:HD11	1:A:477:LEU:HD21	1.86	0.56
1:A:698:ARG:NE	1:A:755:PHE:CE1	2.74	0.55
1:A:608:GLU:OE2	1:A:645:THR:N	2.38	0.54
1:A:675:GLU:HG2	1:A:676:LEU:HD23	1.90	0.53
1:A:676:LEU:CD1	1:A:712:CYS:SG	2.97	0.53
1:B:449:GLY:HA3	1:B:491:LEU:HD23	1.91	0.52
1:A:556:ARG:HA	1:A:559:GLU:HG2	1.92	0.52
1:A:418:ARG:O	1:A:422:ILE:HG12	2.10	0.52
1:A:656:ARG:O	1:A:660:LEU:HG	2.10	0.52
1:A:449:GLY:HA3	1:A:491:LEU:HD23	1.91	0.51
1:A:398:GLU:O	1:A:401:ILE:HG13	2.11	0.51
1:A:712:CYS:HB3	1:A:717:ILE:HB	1.93	0.50
1:B:745:LYS:HE3	1:B:759:VAL:HG22	1.94	0.50
1:A:658:ALA:HB2	1:A:700:LEU:HD21	1.93	0.50
1:A:728:TYR:OH	1:A:738:THR:HG21	2.10	0.50
1:A:676:LEU:HD12	1:A:712:CYS:SG	2.53	0.49
1:A:491:LEU:O	1:A:495:MET:HG2	2.13	0.48
1:A:658:ALA:HB2	1:A:700:LEU:HD11	1.94	0.48
1:A:462:LYS:HD3	1:A:462:LYS:C	2.38	0.48
1:B:565:SER:O	1:B:566:ASP:HB2	2.12	0.48
1:B:557:ILE:HA	1:B:561:LEU:HB2	1.95	0.48
1:B:725:VAL:HG13	1:B:739:PHE:CE1	2.47	0.48
1:A:673:HIS:O	1:A:676:LEU:HG	2.13	0.48
1:A:497:THR:HG23	1:A:546:MET:HE3	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:720:LYS:HB2	1:A:722:LYS:HG2	1.95	0.48
1:A:658:ALA:HB1	1:A:684:PHE:CE2	2.49	0.47
1:A:714:VAL:HG11	1:A:768:ILE:HG22	1.96	0.47
1:A:430:PHE:CD2	1:A:442:GLY:CA	2.97	0.47
1:A:445:ARG:HD2	1:A:495:MET:HE2	1.96	0.47
1:B:475:LYS:O	1:B:476:LEU:C	2.57	0.47
1:A:453:TYR:HE2	1:A:454:TYR:CE1	2.33	0.46
1:A:642:LYS:HG3	1:A:643:LYS:N	2.30	0.46
1:B:516:TRP:CZ2	1:B:520:VAL:HG11	2.50	0.46
1:B:432:LYS:O	1:B:432:LYS:HD3	2.15	0.46
1:A:453:TYR:CE2	1:A:454:TYR:CE1	3.04	0.46
1:A:774:THR:HG1	1:B:511:ASP:N	2.14	0.46
1:A:426:PHE:CD1	1:A:446:TYR:HB2	2.51	0.46
1:A:658:ALA:CA	1:A:700:LEU:HD21	2.46	0.45
1:B:411:PRO:HG3	1:B:477:LEU:HB3	1.97	0.45
1:A:411:PRO:HG3	1:A:477:LEU:HB3	1.99	0.45
1:B:668:ARG:HH11	1:B:668:ARG:CG	2.29	0.45
1:A:672:GLU:O	1:A:672:GLU:HG3	2.15	0.45
1:B:725:VAL:CG1	1:B:739:PHE:CE1	3.00	0.45
1:A:494:VAL:HG23	1:A:495:MET:N	2.32	0.45
1:A:517:ILE:HD11	1:A:521:LEU:HD11	1.99	0.45
1:A:494:VAL:O	1:A:495:MET:C	2.60	0.44
1:B:438:CYS:O	1:B:441:ILE:N	2.46	0.44
1:B:470:ILE:HD12	1:B:470:ILE:H	1.83	0.44
1:B:423:GLY:O	1:B:427:LYS:HG3	2.18	0.44
1:B:491:LEU:O	1:B:495:MET:HG2	2.17	0.44
1:B:404:PHE:N	1:B:404:PHE:CD1	2.86	0.43
1:B:455:ARG:HG2	1:B:455:ARG:HH11	1.83	0.43
1:B:668:ARG:HH11	1:B:668:ARG:HG2	1.83	0.43
1:A:557:ILE:HA	1:A:561:LEU:HB2	2.00	0.43
1:B:445:ARG:NE	1:B:498:TYR:CD1	2.86	0.43
1:B:584:ASP:OD2	1:B:584:ASP:N	2.51	0.43
1:A:698:ARG:HG3	1:A:698:ARG:HH11	1.83	0.43
1:A:457:MET:O	1:A:461:LEU:HG	2.19	0.42
1:B:512:LEU:HD12	1:B:512:LEU:N	2.34	0.42
1:A:399:ASN:O	1:A:402:SER:OG	2.35	0.42
1:A:480:ASN:O	1:A:484:MET:HG2	2.20	0.42
1:B:426:PHE:CD1	1:B:446:TYR:HB2	2.53	0.42
1:B:385:LEU:C	1:B:385:LEU:HD23	2.45	0.42
1:A:758:SER:O	1:A:762:GLN:HG2	2.20	0.42
1:B:438:CYS:HB3	1:B:441:ILE:CD1	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:669:LEU:CD2	1:A:723:ILE:HG22	2.47	0.41
1:A:455:ARG:NH1	1:A:539:GLU:OE2	2.54	0.41
1:A:719:LEU:HD22	1:A:723:ILE:HD12	2.00	0.41
1:A:516:TRP:O	1:A:520:VAL:HG23	2.21	0.41
1:A:656:ARG:NH1	1:A:783:PRO:O	2.53	0.41
1:A:693:GLU:OE2	1:A:763:ARG:NH1	2.54	0.41
1:B:668:ARG:CG	1:B:668:ARG:NH1	2.84	0.41
1:A:516:TRP:CE2	1:A:520:VAL:CG2	3.04	0.41
1:B:392:ALA:O	1:B:451:ARG:NH1	2.41	0.41
1:A:546:MET:HA	1:A:546:MET:HE2	2.02	0.41
1:A:611:ARG:HD3	1:A:611:ARG:HA	1.90	0.41
1:B:430:PHE:CD1	1:B:430:PHE:C	2.98	0.41
1:B:495:MET:HE3	1:B:495:MET:HB3	1.82	0.41
1:B:698:ARG:HA	2:B:840:HOH:O	2.21	0.41
1:A:466:GLU:O	1:A:466:GLU:HG3	2.22	0.40
1:A:516:TRP:CE2	1:A:520:VAL:HG21	2.57	0.40
1:A:747:GLU:O	1:A:747:GLU:CG	2.68	0.40
1:A:513:SER:O	1:A:516:TRP:N	2.46	0.40
1:B:468:LEU:HB3	1:B:472:ASN:HB2	2.04	0.40
1:A:658:ALA:HA	1:A:700:LEU:HD21	2.04	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	326/390 (84%)	321 (98%)	5 (2%)	0	100	100
1	B	345/390 (88%)	340 (99%)	5 (1%)	0	100	100
All	All	671/780 (86%)	661 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	312/362 (86%)	298 (96%)	14 (4%)	24	22
1	B	328/362 (91%)	302 (92%)	26 (8%)	11	7
All	All	640/724 (88%)	600 (94%)	40 (6%)	16	11

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	397	SER
1	A	399	ASN
1	A	421	ASP
1	A	447	LYS
1	A	470	ILE
1	A	471	GLN
1	A	495	MET
1	A	548	LYS
1	A	648	SER
1	A	677	GLU
1	A	698	ARG
1	A	746	GLU
1	A	747	GLU
1	A	766	THR
1	B	383	GLN
1	B	385	LEU
1	B	388	ILE
1	B	393	SER
1	B	399	ASN
1	B	401	ILE
1	B	402	SER
1	B	439	VAL
1	B	440	GLU
1	B	473	PHE
1	B	511	ASP
1	B	520	VAL
1	B	529	TYR

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Mol	Chain	Res	Type
1	B	533	LYS
1	B	575	GLN
1	B	583	THR
1	B	584	ASP
1	B	600	HIS
1	B	608	GLU
1	B	668	ARG
1	B	713	LYS
1	B	714	VAL
1	B	725	VAL
1	B	747	GLU
1	B	748	GLU
1	B	775	ARG

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	405	ASN
1	A	686	HIS
1	B	673	HIS
1	B	685	GLN
1	B	689	GLN
1	B	690	ASN

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	336/390 (86%)	1.17	52 (15%) 5 5	55, 95, 144, 183	0
1	B	353/390 (90%)	0.29	16 (4%) 38 43	23, 48, 92, 122	0
All	All	689/780 (88%)	0.72	68 (9%) 13 14	23, 72, 138, 183	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	612	ALA	4.7
1	B	610	VAL	4.2
1	B	383	GLN	4.0
1	A	724	ILE	3.9
1	B	382	ILE	3.8
1	B	437	GLY	3.8
1	A	761	MET	3.6
1	B	435	GLY	3.6
1	A	743	LEU	3.4
1	A	753	ILE	3.4
1	A	670	LEU	3.4
1	A	719	LEU	3.3
1	B	392	ALA	3.3
1	A	714	VAL	3.2
1	B	439	VAL	3.1
1	A	768	ILE	3.1
1	B	498	TYR	3.1
1	A	735	VAL	3.0
1	A	694	LEU	3.0
1	B	434	VAL	2.8
1	B	438	CYS	2.8
1	A	760	PHE	2.8
1	A	749	TYR	2.8
1	A	782	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	600	HIS	2.7
1	A	723	ILE	2.7
1	A	717	ILE	2.7
1	A	499	SER	2.6
1	A	755	PHE	2.5
1	A	728	TYR	2.5
1	A	758	SER	2.5
1	A	733	HIS	2.5
1	A	711	ILE	2.5
1	A	752	ILE	2.5
1	B	387	MET	2.4
1	B	430	PHE	2.4
1	B	441	ILE	2.4
1	A	756	TYR	2.4
1	A	669	LEU	2.3
1	A	773	SER	2.3
1	A	424	TYR	2.3
1	B	385	LEU	2.3
1	B	541	ASN	2.3
1	B	388	ILE	2.3
1	A	665	LEU	2.3
1	A	772	ALA	2.2
1	A	401	ILE	2.2
1	A	732	PRO	2.2
1	A	779	LEU	2.2
1	A	425	ILE	2.2
1	A	571	ASP	2.2
1	A	731	LEU	2.2
1	A	544	ARG	2.2
1	A	441	ILE	2.1
1	A	575	GLN	2.1
1	A	404	PHE	2.1
1	A	570	PHE	2.1
1	A	610	VAL	2.1
1	A	688	LEU	2.1
1	A	394	ASP	2.1
1	A	388	ILE	2.1
1	A	738	THR	2.1
1	A	766	THR	2.1
1	A	671	SER	2.1
1	A	512	LEU	2.0
1	A	776	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	771	TYR	2.0
1	A	435	GLY	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.