



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:42 PM UTC

PDB ID : 3CQC / pdb_00003cqc
Title : Nucleoporin Nup107/Nup133 interaction complex
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Deposited on : 2008-04-02
Resolution : 2.53 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

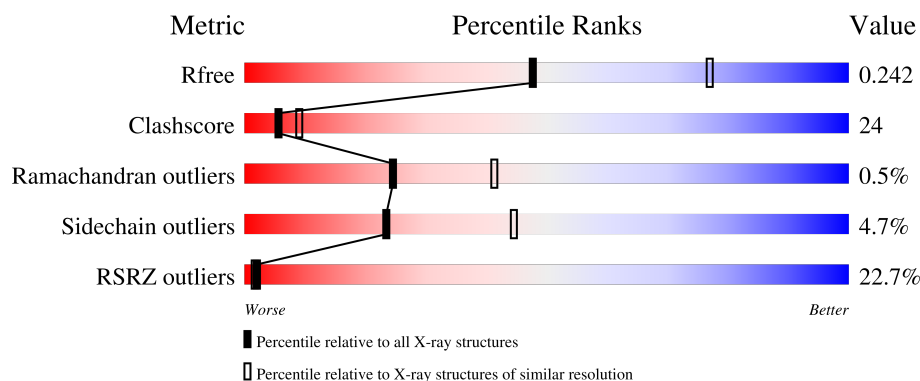
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.53 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	7383 (2.54-2.50)
Clashscore	190562	8079 (2.54-2.50)
Ramachandran outliers	187476	7944 (2.54-2.50)
Sidechain outliers	187428	7946 (2.54-2.50)
RSRZ outliers	180081	7387 (2.54-2.50)

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	270	
2	B	227	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3567 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 Nuclear pore complex protein Nup107.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	0	0
			1993	1282	343	355	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	656	GLY	-	expression tag	UNP P57740
A	657	SER	-	expression tag	UNP P57740

- Molecule 2 is a protein called Nuclear pore complex protein Nup133.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	186	Total	C	N	O	S	76	0	0
			1438	915	238	278	7			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	930	MET	-	expression tag	UNP Q8WUM0
B	931	ALA	-	expression tag	UNP Q8WUM0
B	932	SER	-	expression tag	UNP Q8WUM0
B	933	HIS	-	expression tag	UNP Q8WUM0
B	934	MET	-	expression tag	UNP Q8WUM0

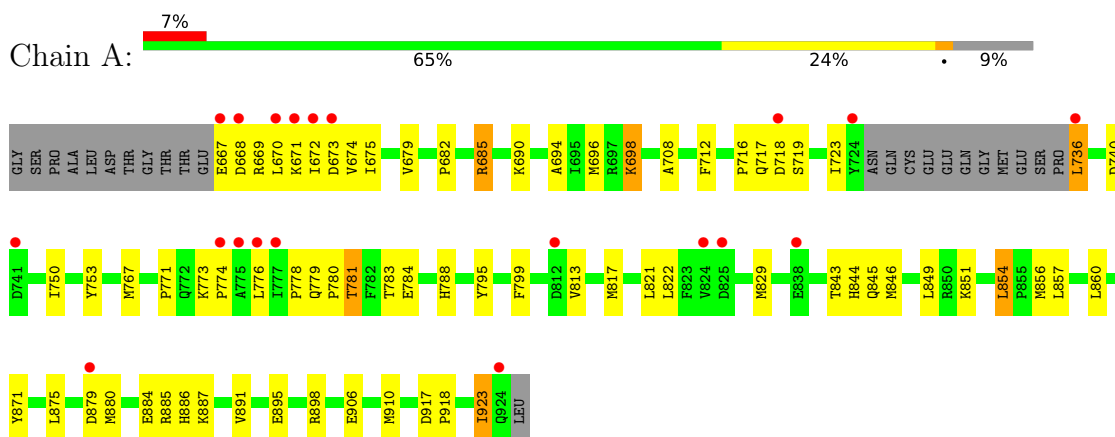
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	114	Total	O	0	0
			114	114		
3	B	22	Total	O	0	0
			22	22		

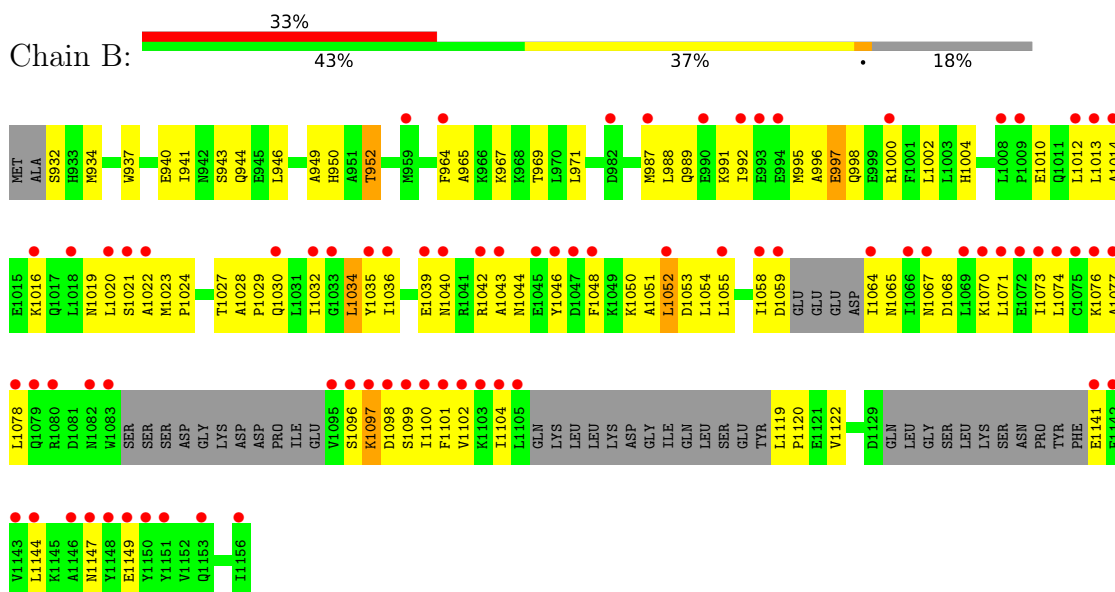
3 Residue-property plots [i](#)

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: Nuclear pore complex protein Nup107



- Molecule 2: Nuclear pore complex protein Nup133



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	51.74Å 127.94Å 152.74Å 90.00° 97.36° 90.00°	Depositor
Resolution (Å)	39.63 – 2.53 39.63 – 2.54	Depositor EDS
% Data completeness (in resolution range)	98.4 (39.63-2.53) 98.4 (39.63-2.54)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.54Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.217 , 0.245 0.216 , 0.242	Depositor DCC
R_{free} test set	1629 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	49.2	Xtriage
Anisotropy	0.358	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3567	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.82% 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	0.65	0/2037	0.86	3/2749 (0.1%)
2	B	0.44	0/1454	0.77	0/1961
All	All	0.57	0/3491	0.82	3/4710 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	854	LEU	CA-C-N	-6.12	112.71	119.19
1	A	854	LEU	C-N-CA	-6.12	112.71	119.19
1	A	781	THR	N-CA-C	-5.23	103.50	110.55

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	1993	0	1981	77	0
2	B	1438	0	1397	84	0
3	A	114	0	0	4	0
3	B	22	0	0	1	0
All	All	3567	0	3378	157	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 24.

All (157) 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)
2:B:1032:ILE:HD11	2:B:1054:LEU:HB2	1.29	1.11
1:A:829:MET:HE3	1:A:846:MET:HG2	1.40	1.02
1:A:696:MET:HE2	1:A:750:ILE:HG21	1.54	0.86
1:A:682:PRO:HA	1:A:685:ARG:HH12	1.40	0.86
2:B:1058:ILE:O	2:B:1059:ASP:HB2	1.72	0.86
1:A:781:THR:HG22	1:A:784:GLU:HG3	1.58	0.85
1:A:698:LYS:HE2	1:A:845:GLN:NE2	1.92	0.84
2:B:1016:LYS:HE2	2:B:1039:GLU:HG3	1.63	0.80
1:A:829:MET:HE2	1:A:849:LEU:HD12	1.62	0.80
2:B:1097:LYS:O	2:B:1101:PHE:HD1	1.65	0.79
2:B:1036:ILE:HG13	2:B:1048:PHE:HE2	1.47	0.78
1:A:696:MET:CE	1:A:712:PHE:HB2	2.13	0.78
1:A:698:LYS:HE2	1:A:845:GLN:HE22	1.50	0.76
1:A:829:MET:CE	1:A:846:MET:HG2	2.16	0.76
1:A:885:ARG:HH11	1:A:885:ARG:HG3	1.49	0.75
2:B:1027:THR:HB	2:B:1029:PRO:HD2	1.70	0.74
2:B:1036:ILE:HG13	2:B:1048:PHE:CE2	2.24	0.72
1:A:696:MET:HE3	1:A:712:PHE:HB2	1.73	0.71
1:A:671:LYS:O	1:A:674:VAL:HG12	1.92	0.70
1:A:844:HIS:HD2	3:A:16:HOH:O	1.75	0.69
2:B:937:TRP:HE3	2:B:941:ILE:HD12	1.58	0.68
1:A:685:ARG:HB2	1:A:685:ARG:HH11	1.58	0.68
2:B:1036:ILE:HD11	2:B:1076:LYS:HB2	1.76	0.68
2:B:1101:PHE:O	2:B:1104:ILE:HG12	1.94	0.67
1:A:696:MET:HE1	1:A:712:PHE:HD1	1.58	0.67
1:A:918:PRO:HD3	2:B:934:MET:HE3	1.77	0.66
2:B:1019:ASN:HB3	2:B:1022:ALA:HB3	1.78	0.66
2:B:1097:LYS:HD2	2:B:1100:ILE:HD11	1.77	0.66
1:A:813:VAL:O	1:A:817:MET:HG2	1.97	0.65
2:B:971:LEU:HD11	2:B:998:GLN:HG2	1.78	0.65
1:A:696:MET:HE1	1:A:712:PHE:CD1	2.32	0.65
1:A:685:ARG:HH21	1:A:716:PRO:HG2	1.62	0.64
2:B:996:ALA:O	2:B:1000:ARG:HG3	1.96	0.64
1:A:696:MET:HE2	1:A:750:ILE:CG2	2.27	0.64
2:B:1077:ALA:HB3	2:B:1097:LYS:HD3	1.79	0.64
2:B:964:PHE:HE1	2:B:1002:LEU:HD22	1.64	0.63
1:A:879:ASP:OD1	2:B:969:THR:HA	1.98	0.63
2:B:937:TRP:CE3	2:B:941:ILE:HD12	2.34	0.63
1:A:781:THR:CG2	1:A:784:GLU:H	2.12	0.62
1:A:781:THR:HG22	1:A:784:GLU:H	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:781:THR:HG22	1:A:784:GLU:CG	2.28	0.62
1:A:671:LYS:HA	1:A:671:LYS:HE2	1.83	0.61
2:B:1020:LEU:H	2:B:1020:LEU:HD22	1.66	0.61
2:B:1097:LYS:HZ3	2:B:1101:PHE:HE1	1.49	0.60
1:A:696:MET:HE1	1:A:712:PHE:HB2	1.83	0.60
2:B:1036:ILE:CD1	2:B:1073:ILE:HA	2.32	0.60
1:A:817:MET:HE1	1:A:860:LEU:HB2	1.83	0.59
2:B:1078:LEU:CD2	2:B:1097:LYS:HG3	2.33	0.59
2:B:1030:GLN:O	2:B:1034:LEU:HB2	2.03	0.58
1:A:781:THR:CG2	1:A:784:GLU:HG3	2.32	0.58
2:B:1073:ILE:HD12	2:B:1074:LEU:N	2.18	0.58
2:B:1052:LEU:HD21	2:B:1073:ILE:HD13	1.85	0.58
1:A:829:MET:HE2	1:A:849:LEU:CD1	2.33	0.57
1:A:885:ARG:HH11	1:A:885:ARG:CG	2.18	0.56
2:B:1065:ASN:HB3	2:B:1068:ASP:HB2	1.87	0.56
1:A:817:MET:HE1	1:A:860:LEU:CB	2.36	0.56
2:B:964:PHE:CE1	2:B:1002:LEU:HD22	2.42	0.55
1:A:767:MET:HE3	3:A:76:HOH:O	2.06	0.55
2:B:1010:GLU:HA	2:B:1013:LEU:CD2	2.36	0.55
1:A:753:TYR:OH	1:A:856:MET:HE1	2.06	0.55
1:A:773:LYS:HD2	1:A:795:TYR:CZ	2.42	0.55
2:B:1004:HIS:CD2	2:B:1054:LEU:HD21	2.42	0.54
1:A:898:ARG:HG2	2:B:946:LEU:HD11	1.89	0.54
1:A:716:PRO:C	1:A:718:ASP:H	2.15	0.53
2:B:988:LEU:O	2:B:992:ILE:HG22	2.09	0.53
2:B:1016:LYS:HE3	2:B:1034:LEU:HD21	1.91	0.53
2:B:1032:ILE:CD1	2:B:1054:LEU:HB2	2.21	0.53
1:A:685:ARG:HB2	1:A:685:ARG:NH1	2.23	0.52
2:B:1012:LEU:HD12	2:B:1012:LEU:C	2.35	0.52
1:A:668:ASP:O	1:A:672:ILE:HG22	2.09	0.52
1:A:669:ARG:HA	1:A:672:ILE:CG2	2.40	0.52
1:A:851:LYS:HA	1:A:891:VAL:HG13	1.91	0.52
1:A:771:PRO:HG2	1:A:799:PHE:HA	1.91	0.52
2:B:943:SER:O	2:B:944:GLN:HB2	2.09	0.51
2:B:1078:LEU:HG	2:B:1097:LYS:HG3	1.92	0.51
2:B:940:GLU:OE1	2:B:952:THR:HG21	2.10	0.51
1:A:773:LYS:HG3	1:A:774:PRO:HD2	1.93	0.51
2:B:1050:LYS:HA	2:B:1053:ASP:OD2	2.10	0.51
1:A:685:ARG:CG	1:A:719:SER:HB3	2.42	0.50
1:A:871:TYR:CZ	1:A:910:MET:HG2	2.47	0.50
2:B:1097:LYS:O	2:B:1100:ILE:HG12	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1013:LEU:C	2:B:1013:LEU:HD12	2.37	0.50
1:A:923:ILE:HD13	2:B:934:MET:HE1	1.93	0.49
2:B:1010:GLU:HA	2:B:1013:LEU:HG	1.95	0.49
1:A:682:PRO:HA	1:A:685:ARG:NH1	2.18	0.49
1:A:669:ARG:O	1:A:673:ASP:HB2	2.12	0.49
1:A:917:ASP:OD1	1:A:917:ASP:C	2.56	0.49
1:A:667:GLU:O	1:A:670:LEU:HB3	2.13	0.49
1:A:696:MET:HE3	1:A:708:ALA:O	2.13	0.48
2:B:1067:ASN:HA	2:B:1070:LYS:CB	2.43	0.48
2:B:1052:LEU:HD21	2:B:1073:ILE:CD1	2.43	0.48
2:B:1099:SER:O	2:B:1102:VAL:HB	2.14	0.48
2:B:971:LEU:HD12	2:B:1002:LEU:HD12	1.95	0.48
1:A:696:MET:CE	1:A:750:ILE:HG21	2.35	0.48
1:A:685:ARG:HE	1:A:716:PRO:HG2	1.79	0.48
1:A:781:THR:HG23	1:A:783:THR:N	2.29	0.47
1:A:685:ARG:HH11	1:A:685:ARG:CB	2.25	0.47
2:B:1052:LEU:HD11	2:B:1073:ILE:CD1	2.46	0.46
1:A:822:LEU:HD21	1:A:880:MET:HG3	1.97	0.46
2:B:1027:THR:CB	2:B:1029:PRO:HD2	2.42	0.46
2:B:1098:ASP:O	2:B:1102:VAL:HG23	2.16	0.46
1:A:690:LYS:HE2	1:A:740:ASP:OD1	2.16	0.46
2:B:1067:ASN:O	2:B:1071:LEU:HG	2.15	0.46
1:A:716:PRO:C	1:A:718:ASP:N	2.74	0.46
1:A:822:LEU:HD22	1:A:886:HIS:CE1	2.51	0.45
1:A:895:GLU:HG3	3:A:12:HOH:O	2.15	0.45
2:B:940:GLU:OE1	2:B:952:THR:CG2	2.64	0.45
2:B:989:GLN:HA	2:B:989:GLN:OE1	2.16	0.45
2:B:1020:LEU:H	2:B:1020:LEU:CD2	2.30	0.45
1:A:884:GLU:OE1	1:A:887:LYS:HE3	2.16	0.45
2:B:1036:ILE:HD12	2:B:1073:ILE:HA	1.99	0.45
2:B:1147:ASN:C	2:B:1149:GLU:H	2.24	0.45
2:B:1052:LEU:HA	2:B:1055:LEU:HD23	1.99	0.44
2:B:940:GLU:OE1	2:B:949:ALA:HA	2.18	0.44
2:B:1032:ILE:HD12	2:B:1051:ALA:O	2.17	0.44
2:B:1048:PHE:O	2:B:1052:LEU:HD22	2.18	0.44
2:B:1036:ILE:HG12	2:B:1036:ILE:O	2.17	0.44
2:B:965:ALA:HB2	2:B:1021:SER:HB2	2.00	0.43
1:A:821:LEU:HD13	1:A:880:MET:HE1	2.00	0.43
1:A:694:ALA:O	1:A:698:LYS:HE3	2.19	0.43
1:A:674:VAL:O	1:A:674:VAL:HG22	2.19	0.43
2:B:1096:SER:O	2:B:1100:ILE:HG23	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:857:LEU:HD23	1:A:857:LEU:HA	1.65	0.43
1:A:906:GLU:O	1:A:910:MET:HE2	2.18	0.43
2:B:1067:ASN:HA	2:B:1070:LYS:HB3	2.01	0.43
1:A:716:PRO:O	1:A:718:ASP:N	2.51	0.43
2:B:1013:LEU:HD12	2:B:1014:ALA:N	2.34	0.43
2:B:1078:LEU:HD11	2:B:1097:LYS:HE3	2.01	0.43
1:A:779:GLN:N	1:A:780:PRO:HD3	2.34	0.43
2:B:950:HIS:HE1	2:B:995:MET:HE1	1.83	0.43
2:B:1028:ALA:O	2:B:1032:ILE:HG12	2.19	0.43
2:B:967:LYS:NZ	3:B:131:HOH:O	2.52	0.42
2:B:1042:ARG:O	2:B:1043:ALA:C	2.60	0.42
2:B:1012:LEU:HD22	2:B:1040:ASN:HB2	2.01	0.42
2:B:987:MET:O	2:B:991:LYS:HG2	2.19	0.42
1:A:885:ARG:HG3	1:A:885:ARG:NH1	2.27	0.42
2:B:1044:ASN:HD21	2:B:1046:TYR:HD2	1.68	0.42
1:A:849:LEU:HD23	1:A:849:LEU:HA	1.84	0.42
1:A:776:LEU:HB3	1:A:778:PRO:HD3	2.02	0.41
2:B:1067:ASN:HA	2:B:1070:LYS:HB2	2.02	0.41
1:A:669:ARG:HA	1:A:669:ARG:HD3	1.68	0.41
2:B:1023:MET:HB2	2:B:1024:PRO:HD2	2.02	0.41
2:B:1019:ASN:O	2:B:1020:LEU:C	2.63	0.41
2:B:1032:ILE:CD1	2:B:1051:ALA:HA	2.50	0.41
2:B:1077:ALA:CB	2:B:1097:LYS:HD3	2.47	0.41
2:B:1012:LEU:HD13	2:B:1034:LEU:HD22	2.02	0.41
2:B:1028:ALA:HB3	2:B:1029:PRO:HD3	2.02	0.41
1:A:719:SER:O	1:A:723:ILE:HG13	2.21	0.41
1:A:776:LEU:HD23	1:A:788:HIS:CE1	2.55	0.41
1:A:854:LEU:HD23	1:A:854:LEU:HA	1.83	0.41
2:B:1073:ILE:CD1	2:B:1074:LEU:HG	2.51	0.41
2:B:1097:LYS:HA	2:B:1100:ILE:HG12	2.02	0.41
1:A:736:LEU:HA	1:A:736:LEU:HD23	1.64	0.41
2:B:997:GLU:OE2	2:B:997:GLU:HA	2.21	0.40
2:B:1141:GLU:HA	2:B:1144:LEU:HB2	2.04	0.40
1:A:685:ARG:HD3	1:A:685:ARG:HA	1.73	0.40
1:A:679:VAL:HA	3:A:65:HOH:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	243/270 (90%)	236 (97%)	6 (2%)	1 (0%)	30	47
2	B	176/227 (78%)	159 (90%)	16 (9%)	1 (1%)	21	36
All	All	419/497 (84%)	395 (94%)	22 (5%)	2 (0%)	24	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	717	GLN
2	B	1120	PRO

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	211/240 (88%)	204 (97%)	7 (3%)	33	58
2	B	148/203 (73%)	138 (93%)	10 (7%)	14	28
All	All	359/443 (81%)	342 (95%)	17 (5%)	23	44

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	675	ILE
1	A	685	ARG
1	A	698	LYS
1	A	736	LEU

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Mol	Chain	Res	Type
1	A	843	THR
1	A	875	LEU
1	A	923	ILE
2	B	932	SER
2	B	952	THR
2	B	997	GLU
2	B	1034	LEU
2	B	1035	TYR
2	B	1052	LEU
2	B	1064	ILE
2	B	1097	LYS
2	B	1119	LEU
2	B	1122	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	768	ASN
1	A	788	HIS
1	A	872	GLN
1	A	886	HIS
2	B	944	GLN
2	B	958	ASN
2	B	998	GLN
2	B	1004	HIS
2	B	1017	GLN
2	B	1040	ASN

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 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	247/270 (91%)	0.53	20 (8%) 18 16	34, 51, 84, 104	0
2	B	175/227 (77%)	2.18	76 (43%) 0 0	45, 114, 163, 167	0
All	All	422/497 (84%)	1.21	96 (22%) 2 2	34, 61, 159, 167	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1105	LEU	10.0
2	B	1104	ILE	8.2
2	B	1150	TYR	7.4
2	B	1071	LEU	7.3
2	B	1066	ILE	7.3
2	B	1064	ILE	7.0
2	B	1147	ASN	6.9
2	B	1142	PHE	6.6
2	B	1069	LEU	6.5
2	B	1020	LEU	5.7
2	B	1083	TRP	5.7
2	B	1074	LEU	5.3
2	B	1075	CYS	5.2
2	B	1036	ILE	5.2
2	B	1102	VAL	5.2
2	B	1021	SER	4.9
2	B	1098	ASP	4.9
2	B	1146	ALA	4.9
1	A	670	LEU	4.7
2	B	1095	VAL	4.5
2	B	1144	LEU	4.5
2	B	1151	TYR	4.3
1	A	667	GLU	4.3
1	A	668	ASP	4.3

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Mol	Chain	Res	Type	RSRZ
2	B	1058	ILE	4.2
2	B	1103	LYS	4.1
2	B	1048	PHE	4.1
2	B	1078	LEU	4.0
2	B	1013	LEU	3.9
2	B	1073	ILE	3.9
2	B	1052	LEU	3.8
2	B	1055	LEU	3.8
2	B	1077	ALA	3.8
2	B	1148	TYR	3.8
2	B	1156	ILE	3.7
1	A	775	ALA	3.7
2	B	1046	TYR	3.7
1	A	924	GLN	3.7
2	B	1143	VAL	3.7
1	A	671	LYS	3.7
1	A	825	ASP	3.7
2	B	1047	ASP	3.7
2	B	1070	LYS	3.5
1	A	776	LEU	3.4
2	B	1097	LYS	3.4
2	B	1141	GLU	3.4
2	B	994	GLU	3.4
2	B	1082	ASN	3.3
2	B	1018	LEU	3.3
2	B	1035	TYR	3.2
2	B	1014	ALA	3.2
2	B	1043	ALA	3.1
2	B	1072	GLU	3.1
2	B	1101	PHE	3.1
2	B	990	GLU	3.0
2	B	1100	ILE	3.0
1	A	718	ASP	3.0
2	B	1099	SER	3.0
1	A	724	TYR	2.9
2	B	1096	SER	2.9
2	B	1009	PRO	2.8
1	A	736	LEU	2.8
1	A	879	ASP	2.8
2	B	1022	ALA	2.8
2	B	1032	ILE	2.7
1	A	672	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	1079	GLN	2.7
1	A	824	VAL	2.7
2	B	987	MET	2.6
2	B	1076	LYS	2.6
2	B	1030	GLN	2.6
2	B	1033	GLY	2.5
1	A	838	GLU	2.5
2	B	1045	GLU	2.5
1	A	673	ASP	2.5
2	B	1059	ASP	2.5
2	B	1008	LEU	2.4
2	B	992	ILE	2.4
2	B	1012	LEU	2.4
1	A	812	ASP	2.3
2	B	1149	GLU	2.3
2	B	1067	ASN	2.3
2	B	1153	GLN	2.3
2	B	1016	LYS	2.3
1	A	741	ASP	2.2
2	B	1000	ARG	2.2
1	A	774	PRO	2.2
2	B	1040	ASN	2.2
2	B	964	PHE	2.2
2	B	993	GLU	2.1
2	B	1042	ARG	2.0
2	B	1080	ARG	2.0
2	B	1039	GLU	2.0
1	A	777	ILE	2.0
2	B	959	MET	2.0
2	B	982	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.