



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:56 AM GMT

PDB ID : 3CQC
Title : Nucleoporin Nup107/Nup133 interaction complex
Authors : Jeudy, S.; Boehmer, T.; Berke, I.; Schwartz, T.U.
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

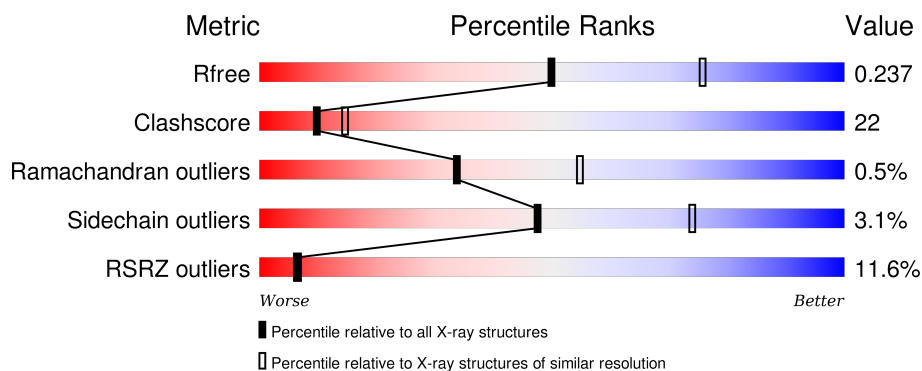
1 Overall quality at a glance

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}	91344	4241 (2.54-2.50)
Clashscore	102246	4968 (2.54-2.50)
Ramachandran outliers	100387	4873 (2.54-2.50)
Sidechain outliers	100360	4875 (2.54-2.50)
RSRZ outliers	91569	4253 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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

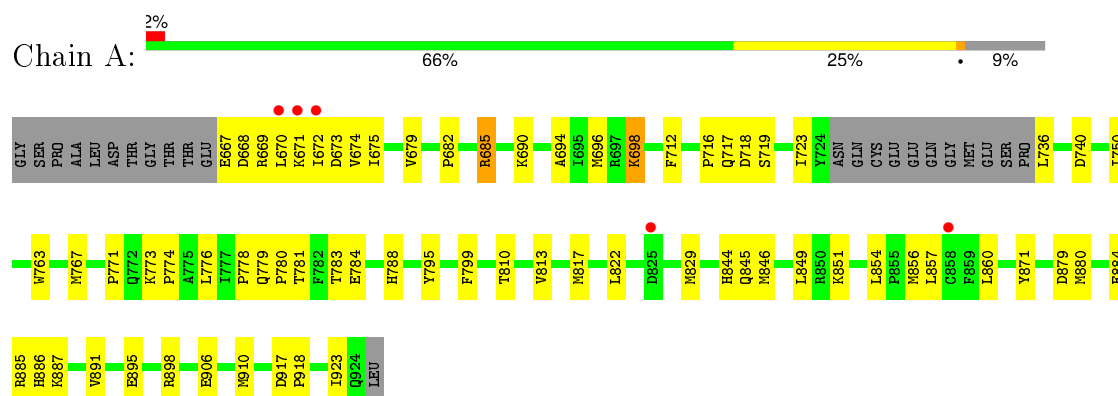
- Molecule 3 is water.

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

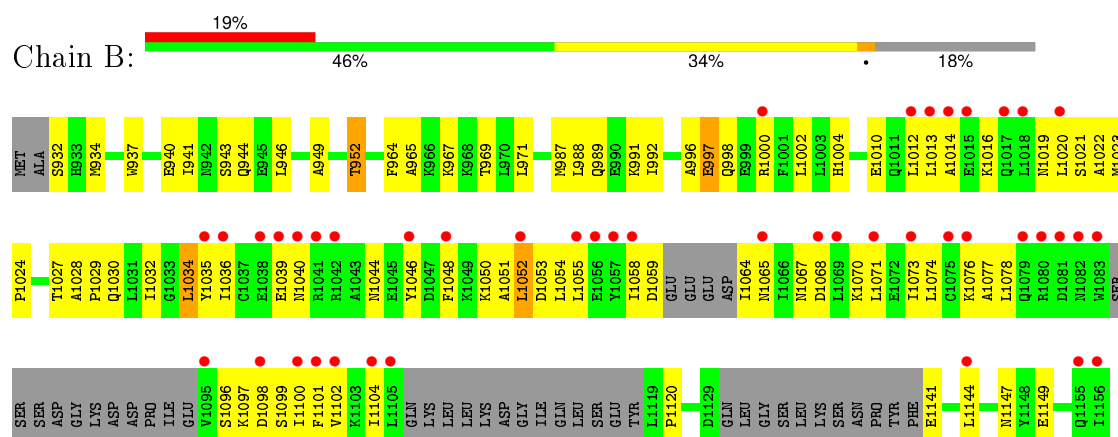
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 – 2.53	Depositor EDS
% Data completeness (in resolution range)	98.4 (39.63-2.53) 98.4 (39.63-2.53)	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 0.210 , 0.237	Depositor DCC
R_{free} test set	1628 reflections (5.06%)	DCC
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 , 57.2	EDS
Estimated twinning fraction	0.027 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 32203 reflections	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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.53	0/2037	0.59	0/2749
2	B	0.35	0/1454	0.46	0/1961
All	All	0.46	0/3491	0.54	0/4710

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 [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	1993	0	1981	71	0
2	B	1438	0	1397	80	0
3	A	114	0	0	4	0
3	B	22	0	0	1	0
All	All	3567	0	3378	148	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 22.

All (148) 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:696:MET:HE3	1:A:712:PHE:HB2	1.25	1.16

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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	1.50	0.93
2:B:1058:ILE:O	2:B:1059:ASP:HB2	1.72	0.86
1:A:682:PRO:HA	1:A:685:ARG:HH12	1.40	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
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: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:696:MET:CE	1:A:712:PHE:HB2	2.13	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: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:767:MET:HE3	3:A:76:HOH:O	1.95	0.67
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
2:B:971:LEU:HD11	2:B:998:GLN:HG2	1.78	0.65
1:A:813:VAL:O	1:A:817:MET:HG2	1.97	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
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
2:B:937:TRP:CE3	2:B:941:ILE:HD12	2.34	0.63
1:A:879:ASP:OD1	2:B:969:THR:HA	1.98	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
1:A:781:THR:HG22	1:A:784:GLU:CG	2.28	0.62
1:A:671:LYS:HE2	1:A:671:LYS:HA	1.83	0.61
2:B:1020:LEU:HD22	2:B:1020:LEU:H	1.66	0.61
1:A:817:MET:HE3	1:A:860:LEU:HB2	1.84	0.60
2:B:1036:ILE:CD1	2:B:1073:ILE:HA	2.32	0.60
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:696:MET:CE	1:A:750:ILE:HG21	2.35	0.57
1:A:885:ARG:CG	1:A:885:ARG:HH11	2.18	0.56
2:B:1065:ASN:HB3	2:B:1068:ASP:HB2	1.87	0.56
2:B:964:PHE:CE1	2:B:1002:LEU:HD22	2.42	0.55
2:B:1010:GLU:HA	2:B:1013:LEU:CD2	2.36	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
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:NH1	1:A:685:ARG:HB2	2.23	0.52
2:B:943:SER:O	2:B:944:GLN:HB2	2.09	0.52
1:A:669:ARG:HA	1:A:672:ILE:CG2	2.40	0.52
1:A:668:ASP:O	1:A:672:ILE:HG22	2.09	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:1078:LEU:HG	2:B:1097:LYS:HG3	1.92	0.51
2:B:1097:LYS:HZ3	2:B:1101:PHE:HE1	1.58	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
2:B:1097:LYS:O	2:B:1100:ILE:HG12	2.12	0.50
1:A:871:TYR:CZ	1:A:910:MET:HG2	2.47	0.50
1:A:696:MET:HE1	1:A:750:ILE:HG21	1.93	0.50
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:667:GLU:O	1:A:670:LEU:HB3	2.13	0.49
1:A:716:PRO:C	1:A:718:ASP:H	2.15	0.49
1:A:763:TRP:HE1	1:A:810:THR:HG1	1.60	0.49
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
1:A:829:MET:HE1	1:A:849:LEU:HD12	1.95	0.48
2:B:971:LEU:HD12	2:B:1002:LEU:HD12	1.95	0.48
1:A:685:ARG:HE	1:A:716:PRO:HG2	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:918:PRO:HD3	2:B:934:MET:HE3	1.95	0.48
1:A:781:THR:HG23	1:A:783:THR:N	2.29	0.47
1:A:685:ARG:CB	1:A:685:ARG:HH11	2.25	0.47
2:B:1012:LEU:HD12	2:B:1012:LEU:C	2.35	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
2:B:1067:ASN:O	2:B:1071:LEU:HG	2.15	0.46
1:A:690:LYS:HE2	1:A:740:ASP:OD1	2.16	0.46
2:B:1020:LEU:CD2	2:B:1020:LEU:H	2.30	0.45
2:B:940:GLU:OE1	2:B:952:THR:CG2	2.64	0.45
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:989:GLN:HA	2:B:989:GLN:OE1	2.16	0.45
1:A:884:GLU:OE1	1:A:887:LYS:HE3	2.16	0.45
2:B:1013:LEU:HD12	2:B:1013:LEU:C	2.37	0.45
2:B:1036:ILE:HD12	2:B:1073:ILE:HA	1.99	0.45
2:B:1052:LEU:HA	2:B:1055:LEU:HD23	1.99	0.44
1:A:917:ASP:C	1:A:917:ASP:OD1	2.56	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:694:ALA:O	1:A:698:LYS:HE3	2.19	0.43
1:A:856:MET:HB3	1:A:856:MET:HE2	1.53	0.43
2:B:1096:SER:O	2:B:1100:ILE:HG23	2.19	0.43
1:A:674:VAL:HG22	1:A:674:VAL:O	2.19	0.43
2:B:1067:ASN:HA	2:B:1070:LYS:HB3	2.01	0.43
1:A:857:LEU:HD23	1:A:857:LEU:HA	1.65	0.43
1:A:716:PRO:O	1:A:718:ASP:N	2.51	0.43
2:B:1078:LEU:HD11	2:B:1097:LYS:HE3	2.01	0.43
2:B:1013:LEU:HD12	2:B:1014:ALA:N	2.34	0.43
2:B:1028:ALA:O	2:B:1032:ILE:HG12	2.19	0.43
1:A:779:GLN:N	1:A:780:PRO:HD3	2.34	0.43
2:B:967:LYS:NZ	3:B:131:HOH:O	2.52	0.42
1:A:696:MET:HE1	1:A:712:PHE:HD1	1.83	0.42
2:B:1012:LEU:HD22	2:B:1040:ASN:HB2	2.01	0.42
1:A:829:MET:CE	1:A:849:LEU:HD12	2.49	0.42
2:B:987:MET:O	2:B:991:LYS:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:885:ARG:NH1	1:A:885:ARG:HG3	2.27	0.42
2:B:1044:ASN:HD21	2:B:1046:TYR:HD2	1.68	0.42
1:A:849:LEU:HA	1:A:849:LEU:HD23	1.84	0.42
2:B:1067:ASN:HA	2:B:1070:LYS:HB2	2.02	0.41
1:A:776:LEU:HB3	1:A:778:PRO:HD3	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: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:1028:ALA:HB3	2:B:1029:PRO:HD3	2.02	0.41
2:B:1012:LEU:HD13	2:B:1034:LEU:HD22	2.02	0.41
1:A:719:SER:O	1:A:723:ILE:HG13	2.21	0.41
1:A:716:PRO:C	1:A:718:ASP:N	2.74	0.41
2:B:1097:LYS:HA	2:B:1100:ILE:HG12	2.02	0.41
2:B:1073:ILE:CD1	2:B:1074:LEU:HG	2.51	0.41
1:A:776:LEU:HD23	1:A:788:HIS:CE1	2.55	0.41
1:A:854:LEU:HA	1:A:854:LEU:HD23	1.83	0.41
1:A:906:GLU:O	1:A:910:MET:HE2	2.21	0.41
1:A:736:LEU:HD23	1:A:736:LEU:HA	1.64	0.41
2:B:997:GLU:HA	2:B:997:GLU:OE2	2.21	0.40
2:B:1141:GLU:HA	2:B:1144:LEU:HB2	2.04	0.40
1:A:685:ARG:HA	1:A:685:ARG:HD3	1.73	0.40
2:B:1147:ASN:C	2:B:1149:GLU:H	2.24	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 [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	243/270 (90%)	236 (97%)	6 (2%)	1 (0%)	39 60

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	176/227 (78%)	159 (90%)	16 (9%)	1 (1%)	30 49
All	All	419/497 (84%)	395 (94%)	22 (5%)	2 (0%)	34 55

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%)	207 (98%)	4 (2%)	65 86
2	B	148/203 (73%)	141 (95%)	7 (5%)	32 55
All	All	359/443 (81%)	348 (97%)	11 (3%)	47 74

All (11) 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	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

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

Mol	Chain	Res	Type
1	A	768	ASN
1	A	886	HIS
1	A	914	GLN
2	B	944	GLN
2	B	958	ASN
2	B	1004	HIS
2	B	1017	GLN

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 carbohydrates 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	247/270 (91%)	0.43	5 (2%) 68 73	34, 51, 84, 104	0
2	B	175/227 (77%)	1.49	44 (25%) 1 1	45, 114, 163, 167	0
All	All	422/497 (84%)	0.87	49 (11%) 6 6	34, 61, 159, 167	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1083	TRP	8.2
2	B	1104	ILE	8.1
2	B	1082	ASN	7.6
2	B	1058	ILE	6.7
2	B	1105	LEU	6.7
2	B	1075	CYS	6.5
2	B	1069	LEU	6.3
2	B	1095	VAL	5.6
1	A	671	LYS	5.4
2	B	1071	LEU	5.3
2	B	1048	PHE	4.9
2	B	1038	GLU	4.8
2	B	1055	LEU	4.7
2	B	1013	LEU	4.6
2	B	1052	LEU	4.6
2	B	1098	ASP	4.5
2	B	1080	ARG	4.4
2	B	1014	ALA	4.3
2	B	1068	ASP	4.1
2	B	1156	ILE	4.0
2	B	1073	ILE	3.8
2	B	1039	GLU	3.7
2	B	1155	GLN	3.6
2	B	1144	LEU	3.5

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Mol	Chain	Res	Type	RSRZ
2	B	1020	LEU	3.5
2	B	1101	PHE	3.2
2	B	1015	GLU	3.0
2	B	1017	GLN	2.9
2	B	1065	ASN	2.9
2	B	1018	LEU	2.9
2	B	1036	ILE	2.8
2	B	1100	ILE	2.7
2	B	1102	VAL	2.7
2	B	1042	ARG	2.6
2	B	1081	ASP	2.6
1	A	672	ILE	2.6
2	B	1076	LYS	2.5
2	B	1057	TYR	2.5
2	B	1056	GLU	2.5
2	B	1012	LEU	2.4
2	B	1040	ASN	2.3
2	B	1041	ARG	2.2
1	A	670	LEU	2.1
2	B	1079	GLN	2.1
1	A	858	CYS	2.1
2	B	1000	ARG	2.1
2	B	1035	TYR	2.1
2	B	1046	TYR	2.1
1	A	825	ASP	2.1

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 carbohydrates 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.