



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:18 PM UTC

PDB ID : 1H2V / pdb_00001h2v
Title : Structure of the human nuclear cap-binding-complex (CBC)
Authors : Mazza, C.; Segref, A.; Mattaj, I.W.; Cusack, S.
Deposited on : 2002-08-16
Resolution : 2.00 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

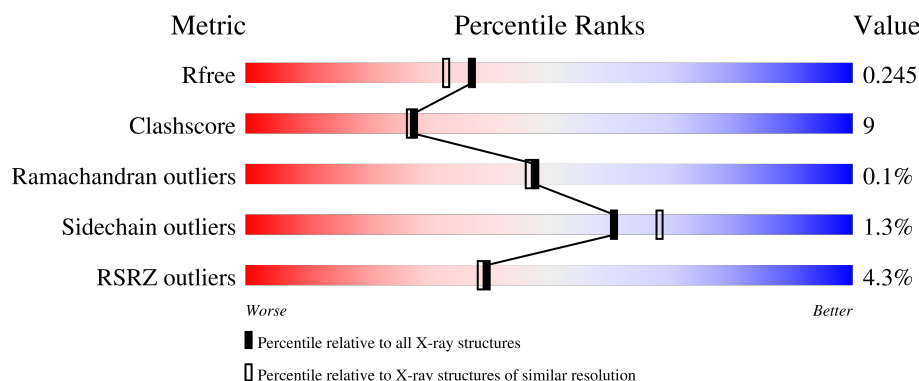
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	C	771	4% 75% 18% 6%
2	Z	156	5% 43% 17% 40%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6961 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 80 KDA NUCLEAR CAP BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	722	Total	C	N	O	S	0	0	0
			5905	3812	996	1059	38			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	479	SER	ALA	engineered mutation	UNP Q09161

- Molecule 2 is a protein called 20 KDA NUCLEAR CAP BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Z	93	Total	C	N	O	S	0	0	0
			762	486	127	143	6			

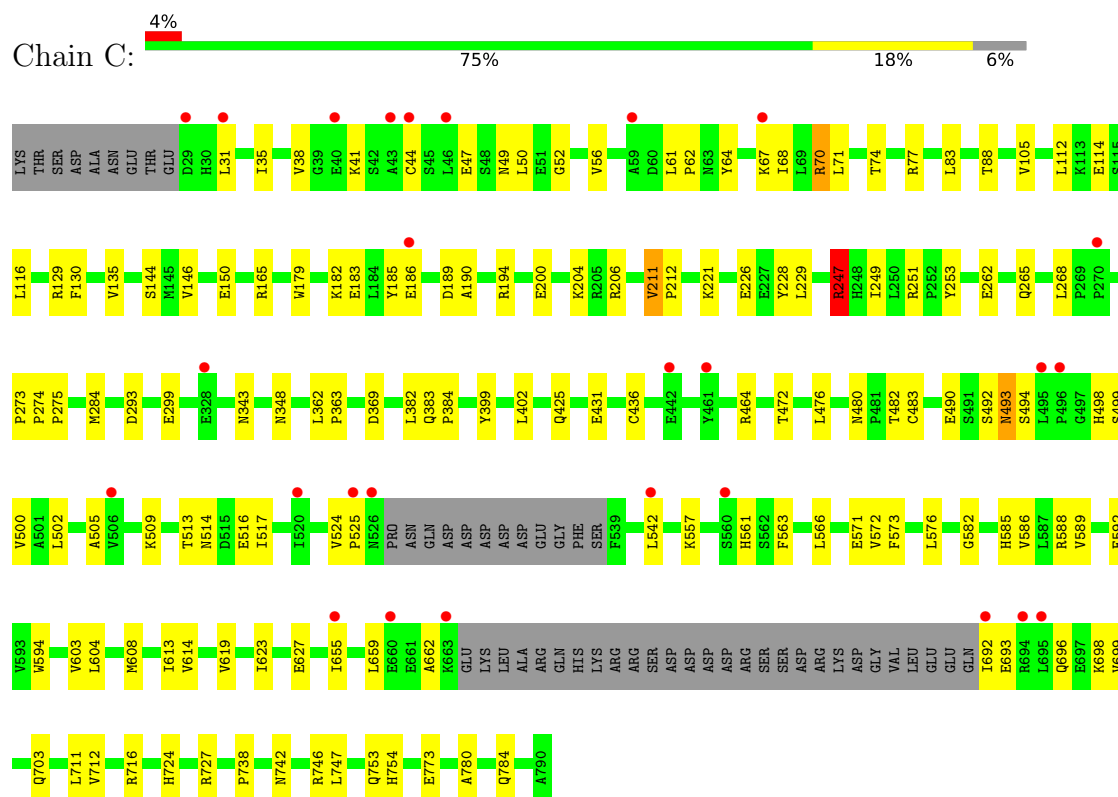
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	250	Total	O	0	0
			250	250		
3	Z	44	Total	O	0	0
			44	44		

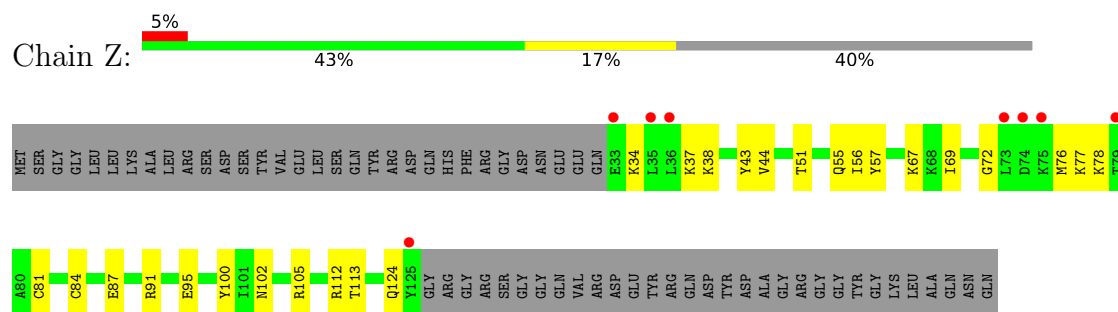
3 Residue-property plots [i](#)

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: 80 KDA NUCLEAR CAP BINDING PROTEIN



• Molecule 2: 20 KDA NUCLEAR CAP BINDING PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	264.14Å 59.60Å 75.43Å 90.00° 99.52° 90.00°	Depositor
Resolution (Å)	19.67 – 2.00 19.67 – 2.00	Depositor EDS
% Data completeness (in resolution range)	79.4 (19.67-2.00) 79.4 (19.67-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.01Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.212 , 0.247 0.211 , 0.245	Depositor DCC
R_{free} test set	3112 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.578	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6961	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% 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	C	0.40	0/6057	0.82	9/8220 (0.1%)
2	Z	0.39	0/775	0.86	3/1035 (0.3%)
All	All	0.39	0/6832	0.83	12/9255 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	247	ARG	N-CA-C	6.18	120.13	112.59
1	C	228	TYR	N-CA-C	6.05	119.49	111.75
1	C	249	ILE	N-CA-C	5.93	117.07	109.30
2	Z	105	ARG	N-CA-C	5.74	119.41	110.17
1	C	165	ARG	N-CA-C	-5.49	104.70	111.40
1	C	226	GLU	N-CA-C	5.44	118.80	110.20
1	C	402	LEU	N-CA-C	5.41	118.68	111.75
2	Z	100	TYR	N-CA-C	5.31	120.77	113.97
1	C	369	ASP	N-CA-C	5.28	117.45	111.11
1	C	211	VAL	CB-CA-C	-5.23	108.73	113.70
2	Z	43	TYR	N-CA-C	-5.14	100.86	109.24
1	C	747	LEU	N-CA-C	-5.01	105.73	111.14

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	C	5905	0	5898	103	0
2	Z	762	0	756	17	0
3	C	250	0	0	9	0
3	Z	44	0	0	3	0
All	All	6961	0	6654	120	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 (120) 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:C:699:VAL:HG12	1:C:703:GLN:HE21	1.30	0.97
1:C:655:ILE:HG22	1:C:699:VAL:HG22	1.63	0.77
2:Z:34:LYS:HD2	2:Z:38:LYS:HZ3	1.50	0.75
1:C:780:ALA:O	1:C:784:GLN:HG3	1.90	0.72
1:C:513:THR:HG22	1:C:516:GLU:HG3	1.72	0.71
1:C:557:LYS:HB3	1:C:561:HIS:HD2	1.58	0.68
1:C:35:ILE:O	1:C:38:VAL:HG12	1.94	0.67
1:C:44:CYS:HB2	1:C:49:ASN:HD21	1.60	0.66
1:C:61:LEU:N	1:C:62:PRO:HD2	2.11	0.65
1:C:588:ARG:HB2	1:C:588:ARG:NH1	2.13	0.64
1:C:182:LYS:O	1:C:186:GLU:HG3	1.98	0.64
1:C:200:GLU:HG2	1:C:204:LYS:HE3	1.79	0.64
1:C:83:LEU:HD11	1:C:130:PHE:HA	1.79	0.63
1:C:608:MET:HE2	1:C:614:VAL:HG13	1.79	0.63
1:C:588:ARG:O	1:C:592:GLU:HG3	1.99	0.63
2:Z:34:LYS:HD2	2:Z:38:LYS:NZ	2.14	0.62
1:C:557:LYS:HB3	1:C:561:HIS:CD2	2.38	0.59
1:C:262:GLU:O	1:C:265:GLN:HG3	2.04	0.58
1:C:582:GLY:O	1:C:586:VAL:HG23	2.03	0.58
2:Z:44:VAL:HG22	2:Z:113:THR:HG22	1.86	0.58
1:C:146:VAL:O	1:C:150:GLU:HG3	2.04	0.57
1:C:67:LYS:HG2	1:C:70:ARG:HH21	1.68	0.57
2:Z:112:ARG:HG2	3:Z:2029:HOH:O	2.04	0.57
1:C:588:ARG:HB2	1:C:588:ARG:HH11	1.67	0.57
1:C:382:LEU:C	1:C:384:PRO:HD3	2.29	0.56
1:C:135:VAL:HG12	1:C:183:GLU:HG2	1.86	0.56
1:C:493:ASN:N	1:C:493:ASN:HD22	2.02	0.56
2:Z:77:LYS:O	2:Z:78:LYS:HB2	2.06	0.56
2:Z:56:ILE:HG22	2:Z:69:ILE:HD13	1.88	0.55
1:C:699:VAL:HG12	1:C:703:GLN:NE2	2.13	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:247:ARG:HB3	1:C:343:ASN:ND2	2.23	0.54
1:C:431:GLU:H	1:C:431:GLU:CD	2.16	0.53
1:C:105:VAL:CG1	1:C:268:LEU:HD23	2.39	0.53
1:C:425:GLN:HG3	3:C:2160:HOH:O	2.08	0.53
1:C:753:GLN:HG2	1:C:754:HIS:CE1	2.44	0.53
1:C:716:ARG:HA	1:C:716:ARG:HH11	1.73	0.52
1:C:586:VAL:CG1	1:C:608:MET:HE1	2.39	0.52
2:Z:102:ASN:HB2	2:Z:113:THR:OG1	2.10	0.52
1:C:247:ARG:HB3	1:C:343:ASN:HD22	1.75	0.52
1:C:492:SER:C	1:C:494:SER:H	2.18	0.52
1:C:383:GLN:N	1:C:384:PRO:HD3	2.25	0.51
1:C:480:ASN:ND2	1:C:482:THR:HG23	2.25	0.51
1:C:573:PHE:HB3	1:C:613:ILE:HG23	1.90	0.51
1:C:524:VAL:HG13	1:C:542:LEU:HD23	1.93	0.51
1:C:47:GLU:H	1:C:47:GLU:CD	2.18	0.51
1:C:67:LYS:HZ3	1:C:71:LEU:HD11	1.76	0.51
1:C:557:LYS:HD3	1:C:561:HIS:NE2	2.26	0.50
2:Z:67:LYS:HD3	2:Z:87:GLU:OE2	2.11	0.50
1:C:52:GLY:O	1:C:56:VAL:HG23	2.11	0.50
1:C:284:MET:HE1	1:C:363:PRO:HG3	1.95	0.49
1:C:41:LYS:HD2	3:C:2002:HOH:O	2.13	0.49
1:C:500:VAL:HG11	1:C:524:VAL:HG22	1.94	0.49
1:C:572:VAL:O	1:C:576:LEU:HD13	2.12	0.49
1:C:712:VAL:O	1:C:716:ARG:HG2	2.12	0.49
1:C:592:GLU:HG2	3:C:2206:HOH:O	2.13	0.48
1:C:724:HIS:ND1	1:C:727:ARG:NH2	2.61	0.48
1:C:129:ARG:HD2	3:C:2019:HOH:O	2.12	0.48
1:C:499:SER:HA	1:C:502:LEU:HD12	1.94	0.48
1:C:31:LEU:HD12	1:C:31:LEU:H	1.79	0.47
1:C:74:THR:HA	1:C:77:ARG:NH1	2.29	0.47
1:C:492:SER:C	1:C:494:SER:N	2.72	0.47
2:Z:37:LYS:C	2:Z:38:LYS:HD2	2.38	0.47
1:C:586:VAL:HG11	1:C:608:MET:HE1	1.97	0.47
2:Z:38:LYS:HD2	2:Z:38:LYS:N	2.30	0.47
1:C:498:HIS:O	1:C:502:LEU:HG	2.15	0.47
1:C:513:THR:CG2	1:C:516:GLU:HG3	2.43	0.47
2:Z:57:TYR:CD2	2:Z:69:ILE:HD12	2.49	0.46
1:C:251:ARG:HB3	1:C:253:TYR:CE1	2.50	0.46
1:C:493:ASN:HA	1:C:498:HIS:CG	2.50	0.46
1:C:588:ARG:HD3	1:C:627:GLU:CD	2.40	0.46
1:C:35:ILE:HD12	1:C:71:LEU:HD13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:480:ASN:HD22	1:C:482:THR:HG23	1.81	0.46
1:C:493:ASN:N	1:C:493:ASN:ND2	2.64	0.46
2:Z:56:ILE:HD13	2:Z:84:CYS:SG	2.56	0.46
1:C:472:THR:HG23	3:C:2187:HOH:O	2.15	0.46
1:C:738:PRO:HG2	3:C:2168:HOH:O	2.15	0.46
1:C:200:GLU:CG	1:C:204:LYS:HE3	2.46	0.45
1:C:493:ASN:HA	1:C:498:HIS:CD2	2.52	0.45
1:C:70:ARG:HE	1:C:70:ARG:HB3	1.56	0.45
1:C:514:ASN:OD1	1:C:571:GLU:HB2	2.17	0.45
1:C:31:LEU:HD12	1:C:31:LEU:N	2.32	0.45
1:C:293:ASP:C	1:C:293:ASP:OD2	2.59	0.45
2:Z:78:LYS:HE3	3:Z:2022:HOH:O	2.16	0.44
2:Z:78:LYS:HG2	3:Z:2022:HOH:O	2.17	0.44
2:Z:91:ARG:O	2:Z:95:GLU:HG3	2.16	0.44
1:C:500:VAL:HG11	1:C:524:VAL:CG2	2.47	0.44
1:C:190:ALA:O	1:C:194:ARG:HG3	2.18	0.44
1:C:505:ALA:O	1:C:509:LYS:HG3	2.16	0.44
1:C:483:CYS:HB2	1:C:594:TRP:CH2	2.52	0.43
1:C:742:ASN:O	1:C:746:ARG:HG2	2.18	0.43
1:C:493:ASN:HD22	1:C:494:SER:N	2.17	0.43
1:C:655:ILE:HD13	1:C:698:LYS:HG2	2.01	0.43
2:Z:72:GLY:HA2	2:Z:124:GLN:O	2.19	0.43
1:C:221:LYS:HE3	1:C:221:LYS:HB2	1.89	0.43
1:C:284:MET:HE1	1:C:363:PRO:CG	2.49	0.43
1:C:61:LEU:N	1:C:62:PRO:CD	2.81	0.43
1:C:129:ARG:HD3	1:C:179:TRP:CZ3	2.53	0.43
1:C:362:LEU:HA	1:C:363:PRO:C	2.44	0.43
1:C:348:ASN:HB2	3:C:2125:HOH:O	2.19	0.43
1:C:566:LEU:HD11	1:C:604:LEU:HD22	2.00	0.43
1:C:273:PRO:HA	1:C:274:PRO:HD3	1.92	0.43
1:C:50:LEU:HD13	1:C:88:THR:HG21	2.00	0.42
1:C:144:SER:HB3	1:C:268:LEU:HD12	2.01	0.42
1:C:619:VAL:O	1:C:623:ILE:HG12	2.20	0.42
2:Z:51:THR:HA	2:Z:55:GLN:OE1	2.20	0.42
1:C:185:TYR:O	1:C:189:ASP:HB3	2.20	0.42
1:C:513:THR:O	1:C:517:ILE:HD13	2.20	0.42
1:C:525:PRO:HG3	3:C:2198:HOH:O	2.19	0.42
1:C:662:ALA:O	1:C:692:ILE:HD13	2.20	0.42
1:C:211:VAL:HB	1:C:212:PRO:HD3	2.02	0.42
1:C:476:LEU:HD12	1:C:476:LEU:N	2.35	0.41
1:C:659:LEU:HD21	1:C:696:GLN:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:585:HIS:O	1:C:589:VAL:HG23	2.20	0.41
1:C:692:ILE:HG23	1:C:693:GLU:N	2.35	0.41
1:C:399:TYR:OH	1:C:436:CYS:HB3	2.21	0.41
1:C:563:PHE:HZ	1:C:603:VAL:HB	1.86	0.40
1:C:112:LEU:O	1:C:116:LEU:HG	2.21	0.40
1:C:275:PRO:HD2	3:C:2081:HOH:O	2.21	0.40
1:C:64:TYR:O	1:C:68:ILE:HG13	2.22	0.40
1:C:206:ARG:CZ	1:C:229:LEU:HD12	2.51	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	C	716/771 (93%)	691 (96%)	24 (3%)	1 (0%)	48	46
2	Z	91/156 (58%)	89 (98%)	2 (2%)	0	100	100
All	All	807/927 (87%)	780 (97%)	26 (3%)	1 (0%)	48	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	490	GLU

5.3.2 Protein sidechains [i](#)

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	C	663/708 (94%)	655 (99%)	8 (1%)	63	70
2	Z	81/130 (62%)	79 (98%)	2 (2%)	42	45
All	All	744/838 (89%)	734 (99%)	10 (1%)	61	68

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	70	ARG
1	C	114	GLU
1	C	247	ARG
1	C	299	GLU
1	C	464	ARG
1	C	493	ASN
1	C	711	LEU
1	C	773	GLU
2	Z	76	MET
2	Z	81	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	49	ASN
1	C	63	ASN
1	C	111	GLN
1	C	157	GLN
1	C	198	ASN
1	C	215	GLN
1	C	325	HIS
1	C	343	ASN
1	C	367	HIS
1	C	406	ASN
1	C	439	GLN
1	C	480	ASN
1	C	493	ASN
1	C	553	HIS
1	C	561	HIS
1	C	585	HIS
1	C	596	ASN
1	C	649	ASN
1	C	656	GLN
1	C	696	GLN

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Mol	Chain	Res	Type
1	C	703	GLN
1	C	706	GLN
1	C	715	GLN
1	C	753	GLN
1	C	754	HIS
1	C	756	GLN
1	C	760	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 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	C	722/771 (93%)	0.27	27 (3%) 45 44	27, 47, 90, 110	0
2	Z	93/156 (59%)	0.36	8 (8%) 16 15	30, 46, 84, 92	0
All	All	815/927 (87%)	0.28	35 (4%) 40 39	27, 47, 89, 110	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	461	TYR	3.6
1	C	692	ILE	3.5
1	C	663	LYS	3.4
1	C	526	ASN	3.3
2	Z	35	LEU	3.1
2	Z	36	LEU	3.1
1	C	59	ALA	3.0
2	Z	125	TYR	3.0
2	Z	73	LEU	2.8
1	C	186	GLU	2.8
1	C	660	GLU	2.8
1	C	506	VAL	2.8
1	C	560	SER	2.7
1	C	29	ASP	2.6
1	C	495	LEU	2.6
1	C	270	PRO	2.6
1	C	520	ILE	2.5
2	Z	79	THR	2.4
1	C	542	LEU	2.4
2	Z	33	GLU	2.4
1	C	328	GLU	2.4
1	C	655	ILE	2.4
1	C	43	ALA	2.3
1	C	695	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	40	GLU	2.3
2	Z	74	ASP	2.3
1	C	31	LEU	2.3
1	C	46	LEU	2.3
1	C	496	PRO	2.2
1	C	44	CYS	2.1
1	C	525	PRO	2.1
1	C	67	LYS	2.1
1	C	442	GLU	2.0
2	Z	75	LYS	2.0
1	C	694	ARG	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.