



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:54 AM UTC

PDB ID : 5E0T / pdb_00005e0t
Title : Human PCNA mutant - S228I
Authors : Duffy, C.M.; Hilbert, B.J.; Kelch, B.A.
Deposited on : 2015-09-29
Resolution : 2.67 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

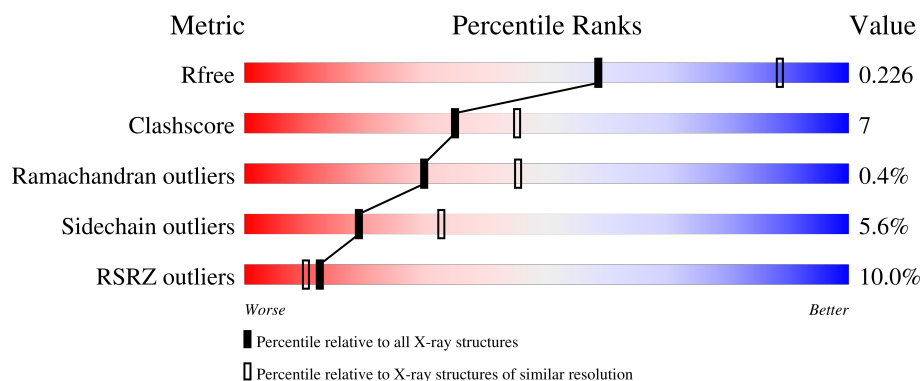
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.67 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	261	10% 75% 18% • 5%
1	B	261	11% 75% 19% • • •
1	C	261	8% 81% 14% • •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6040 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
1	A	248	Total	C	N	O	S	0	0	0
			1909	1204	313	376	16			
1	B	253	Total	C	N	O	S	0	0	0
			1953	1228	318	391	16			
1	C	252	Total	C	N	O	S	0	0	0
			1942	1222	317	387	16			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	ILE	SER	engineered mutation	UNP P12004
B	228	ILE	SER	engineered mutation	UNP P12004
C	228	ILE	SER	engineered mutation	UNP P12004

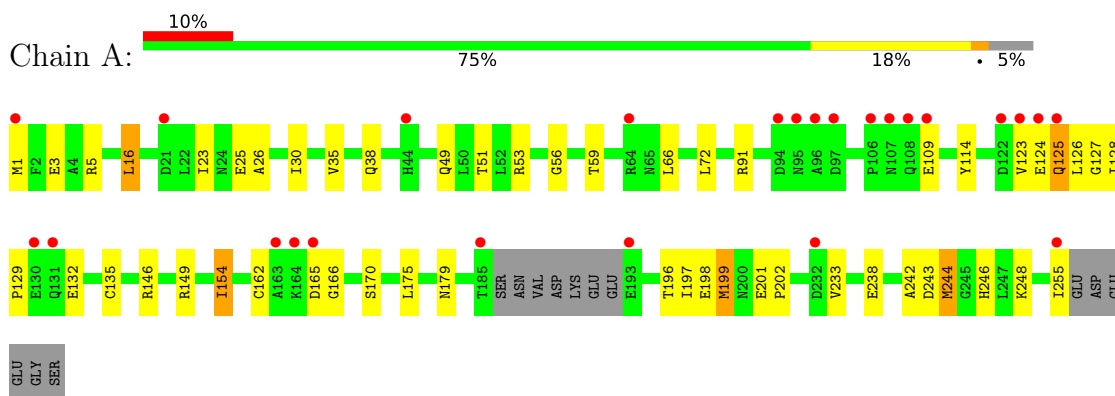
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	70	Total	O	0	0
			70	70		
2	B	74	Total	O	0	0
			74	74		
2	C	92	Total	O	0	0
			92	92		

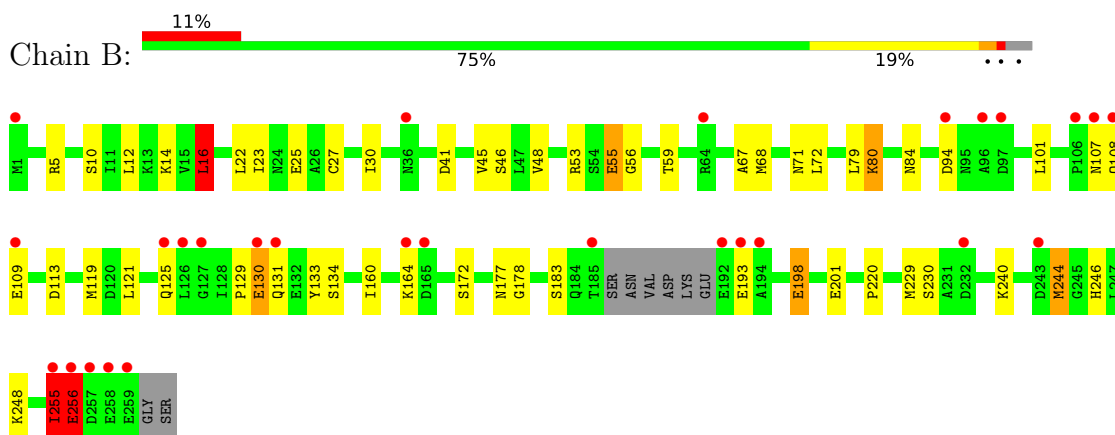
3 Residue-property plots

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.

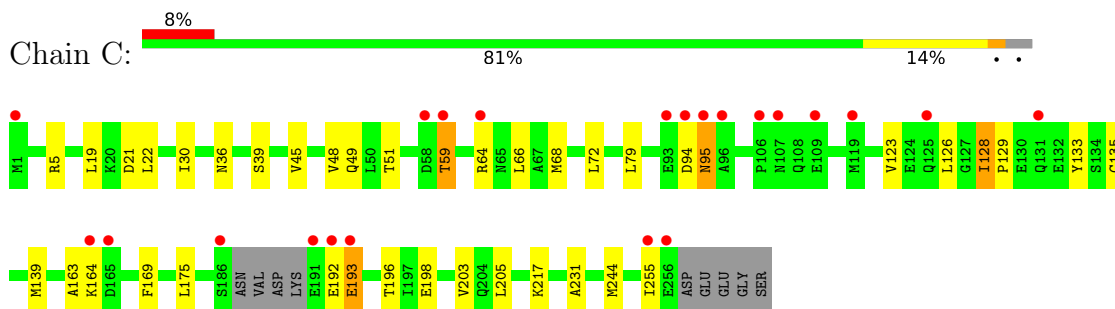
- Molecule 1: Proliferating cell nuclear antigen



- Molecule 1: Proliferating cell nuclear antigen



- Molecule 1: Proliferating cell nuclear antigen



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	162.32Å 162.32Å 139.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.85 – 2.67 42.85 – 2.67	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.85-2.67) 99.7 (42.85-2.67)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.189 , 0.226 0.192 , 0.226	Depositor DCC
R_{free} test set	2143 reflections (3.99%)	wwPDB-VP
Wilson B-factor (Å ²)	47.6	Xtriage
Anisotropy	0.207	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6040	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.23% 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	0.81	0/1934	1.06	7/2612 (0.3%)
1	B	0.83	0/1978	1.07	7/2671 (0.3%)
1	C	0.88	0/1967	1.08	8/2656 (0.3%)
All	All	0.84	0/5879	1.07	22/7939 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	133	TYR	N-CA-C	6.55	120.45	111.54
1	B	108	GLN	N-CA-C	6.53	120.25	112.93
1	C	95	ASN	CA-C-N	-6.23	113.55	122.77
1	C	95	ASN	C-N-CA	-6.23	113.55	122.77
1	C	45	VAL	N-CA-C	-6.19	106.73	111.62
1	A	16	LEU	CA-CB-CG	-6.11	94.93	116.30
1	B	244	MET	N-CA-C	6.02	119.81	112.23
1	B	94	ASP	N-CA-C	5.82	117.70	111.36
1	B	16	LEU	CA-CB-CG	-5.76	96.15	116.30
1	A	198	GLU	CA-C-N	-5.58	115.19	123.11
1	A	198	GLU	C-N-CA	-5.58	115.19	123.11
1	B	80	LYS	N-CA-C	-5.31	106.76	113.18
1	A	244	MET	N-CA-C	5.27	118.97	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	84	ASN	N-CA-C	5.22	117.71	111.71
1	C	126	LEU	CA-C-N	-5.21	115.93	121.45
1	C	126	LEU	C-N-CA	-5.21	115.93	121.45
1	A	154	ILE	N-CA-C	5.13	117.02	111.58
1	A	132	GLU	N-CA-C	5.11	119.79	113.50
1	B	133	TYR	N-CA-C	5.10	118.48	111.54
1	C	198	GLU	CA-C-N	-5.06	115.93	123.11
1	C	198	GLU	C-N-CA	-5.06	115.93	123.11
1	A	242	ALA	CB-CA-C	-5.01	110.78	116.54

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	124	GLU	Peptide
1	B	107	ASN	Peptide

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	1909	0	1929	29	0
1	B	1953	0	1957	41	1
1	C	1942	0	1952	19	0
2	A	70	0	0	1	0
2	B	74	0	0	5	0
2	C	92	0	0	2	0
All	All	6040	0	5838	87	1

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 7.

All (87) 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:3:GLU:OE2	1:A:91:ARG:NH1	2.08	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:139:MET:HE1	1:C:169:PHE:HZ	1.54	0.72
1:B:53:ARG:HD2	2:B:304:HOH:O	1.91	0.69
1:B:255:ILE:HD12	1:B:255:ILE:H	1.56	0.68
1:C:164:LYS:O	2:C:301:HOH:O	2.13	0.67
1:B:41:ASP:OD1	1:B:46:SER:HB3	1.97	0.65
1:A:49:GLN:HG3	1:A:128:ILE:HD11	1.79	0.65
1:C:139:MET:HE1	1:C:169:PHE:CZ	2.34	0.63
1:B:130:GLU:H	1:B:130:GLU:CD	2.07	0.62
1:C:36:ASN:HD21	1:C:49:GLN:HE21	1.46	0.62
1:B:45:VAL:HG22	2:B:301:HOH:O	2.00	0.61
1:C:30:ILE:HG13	1:C:68:MET:HE2	1.80	0.61
1:C:64:ARG:NH2	1:C:94:ASP:O	2.33	0.61
1:C:5:ARG:HB3	1:C:59:THR:HB	1.87	0.56
1:A:5:ARG:HB3	1:A:59:THR:HB	1.85	0.56
1:A:199:MET:HE3	1:A:201:GLU:C	2.30	0.56
1:A:125:GLN:NE2	1:A:126:LEU:H	2.03	0.55
1:A:238:GLU:OE2	1:A:248:LYS:HE2	2.07	0.55
1:B:256:GLU:OE1	1:B:256:GLU:N	2.40	0.55
1:A:199:MET:HE3	1:A:202:PRO:N	2.23	0.54
1:A:38:GLN:HG2	1:A:49:GLN:HG2	1.89	0.54
1:B:255:ILE:HG22	1:B:256:GLU:OE1	2.08	0.54
1:B:5:ARG:HB3	1:B:59:THR:HB	1.90	0.53
1:B:255:ILE:HD12	1:B:255:ILE:N	2.23	0.53
1:B:130:GLU:OE1	1:B:130:GLU:N	2.30	0.53
1:A:246:HIS:CD2	1:A:248:LYS:HG3	2.44	0.52
1:A:166:GLY:HA2	1:A:197:ILE:HD13	1.92	0.51
1:B:14:LYS:HD3	1:B:220:PRO:HB2	1.93	0.51
1:B:240:LYS:NZ	2:B:302:HOH:O	2.45	0.50
1:B:30:ILE:HD11	1:B:68:MET:HE2	1.93	0.50
1:B:25:GLU:HB3	1:B:121:LEU:HD11	1.93	0.50
1:B:27:CYS:SG	1:B:67:ALA:HB1	2.52	0.49
1:A:30:ILE:HB	1:A:66:LEU:HB2	1.95	0.49
1:A:53:ARG:NH2	1:A:243:ASP:O	2.45	0.49
1:B:130:GLU:CD	1:B:130:GLU:N	2.71	0.48
1:B:10:SER:O	1:B:14:LYS:HG3	2.13	0.48
1:B:255:ILE:O	1:B:256:GLU:HB3	2.14	0.47
1:A:53:ARG:NH1	2:A:303:HOH:O	2.48	0.47
1:A:125:GLN:NE2	1:A:127:GLY:H	2.13	0.47
1:C:30:ILE:HB	1:C:66:LEU:HB2	1.96	0.47
1:B:109:GLU:H	1:B:109:GLU:CD	2.24	0.46
1:B:71:ASN:HB2	1:B:119:MET:HE3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:SER:CB	1:B:177:ASN:HB3	2.46	0.46
1:A:109:GLU:OE1	1:B:183:SER:HB2	2.16	0.45
1:C:19:LEU:CD2	1:C:48:VAL:HG11	2.46	0.45
1:B:240:LYS:NZ	2:B:303:HOH:O	2.27	0.45
1:B:240:LYS:HA	1:B:240:LYS:HD2	1.71	0.45
1:A:244:MET:HE2	1:A:244:MET:HB3	1.86	0.45
1:C:30:ILE:CG1	1:C:68:MET:HE2	2.47	0.45
1:A:125:GLN:CD	1:A:126:LEU:H	2.25	0.45
1:C:163:ALA:HB1	2:C:347:HOH:O	2.17	0.45
1:B:53:ARG:NH2	2:B:302:HOH:O	2.24	0.44
1:A:114:TYR:CD1	1:B:178:GLY:HA3	2.53	0.44
1:A:1:MET:HE3	1:A:3:GLU:HB2	2.00	0.44
1:C:193:GLU:H	1:C:193:GLU:HG2	1.51	0.43
1:B:244:MET:HE2	1:B:244:MET:HB3	1.70	0.43
1:B:16:LEU:HD13	1:B:79:LEU:CD1	2.49	0.43
1:B:23:ILE:HG13	1:B:72:LEU:CD1	2.49	0.43
1:B:55:GLU:H	1:B:55:GLU:CD	2.26	0.43
1:C:205:LEU:HD21	1:C:231:ALA:HA	2.02	0.42
1:B:12:LEU:HD12	1:B:12:LEU:HA	1.78	0.42
1:B:30:ILE:CD1	1:B:68:MET:HE2	2.48	0.42
1:B:229:MET:HE2	1:B:229:MET:HB2	1.71	0.42
1:A:3:GLU:CD	1:A:91:ARG:NH1	2.76	0.42
1:B:160:ILE:HG21	1:B:229:MET:HE1	2.02	0.42
1:C:19:LEU:HD21	1:C:48:VAL:HG11	2.02	0.42
1:B:23:ILE:HG22	1:B:41:ASP:HA	2.02	0.42
1:A:146:ARG:HE	1:A:149:ARG:HH22	1.66	0.42
1:A:135:CYS:SG	1:A:162:CYS:HB3	2.60	0.41
1:B:22:LEU:HD23	1:B:48:VAL:HG21	2.01	0.41
1:B:101:LEU:HD12	1:B:101:LEU:N	2.35	0.41
1:B:134:SER:HB2	1:B:230:SER:HA	2.03	0.41
1:C:64:ARG:NE	1:C:94:ASP:HB3	2.36	0.41
1:C:128:ILE:HA	1:C:129:PRO:HD3	1.92	0.41
1:A:30:ILE:HD12	1:A:35:VAL:HG22	2.03	0.41
1:A:56:GLY:HA3	1:A:244:MET:HG2	2.02	0.41
1:A:23:ILE:HG21	1:A:26:ALA:HB2	2.02	0.41
1:B:129:PRO:HA	1:B:130:GLU:OE1	2.21	0.41
1:C:255:ILE:HD13	1:C:255:ILE:HG21	1.85	0.41
1:A:154:ILE:HD13	1:A:175:LEU:HD11	2.02	0.40
1:B:56:GLY:HA3	1:B:244:MET:HE3	2.02	0.40
1:C:21:ASP:OD2	1:C:217:LYS:NZ	2.47	0.40
1:A:23:ILE:HB	1:A:72:LEU:HD12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:ILE:HA	1:A:129:PRO:HD3	1.88	0.40
1:B:246:HIS:CD2	1:B:248:LYS:HG3	2.56	0.40
1:A:170:SER:OG	1:A:179:ASN:HB3	2.21	0.40
1:C:22:LEU:HD12	1:C:22:LEU:HA	1.90	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:GLU:OE2	1:B:198:GLU:OE2[8_555]	1.89	0.31

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	244/261 (94%)	241 (99%)	2 (1%)	1 (0%)	30	45
1	B	249/261 (95%)	245 (98%)	2 (1%)	2 (1%)	16	27
1	C	248/261 (95%)	245 (99%)	3 (1%)	0	100	100
All	All	741/783 (95%)	731 (99%)	7 (1%)	3 (0%)	30	45

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	256	GLU
1	B	255	ILE
1	A	165	ASP

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	216/228 (95%)	207 (96%)	9 (4%)	26	45
1	B	221/228 (97%)	208 (94%)	13 (6%)	18	30
1	C	220/228 (96%)	205 (93%)	15 (7%)	14	25
All	All	657/684 (96%)	620 (94%)	37 (6%)	19	32

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	25	GLU
1	A	51	THR
1	A	123	VAL
1	A	125	GLN
1	A	196	THR
1	A	199	MET
1	A	233	VAL
1	A	255	ILE
1	B	16	LEU
1	B	55	GLU
1	B	80	LYS
1	B	113	ASP
1	B	125	GLN
1	B	130	GLU
1	B	131	GLN
1	B	164	LYS
1	B	193	GLU
1	B	198	GLU
1	B	201	GLU
1	B	255	ILE
1	B	256	GLU
1	C	39	SER
1	C	51	THR
1	C	59	THR
1	C	72	LEU

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Mol	Chain	Res	Type
1	C	79	LEU
1	C	95	ASN
1	C	123	VAL
1	C	128	ILE
1	C	135	CYS
1	C	175	LEU
1	C	192	GLU
1	C	193	GLU
1	C	196	THR
1	C	203	VAL
1	C	244	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	GLN
1	A	213	ASN
1	B	65	ASN
1	B	177	ASN
1	B	179	ASN
1	C	49	GLN
1	C	84	ASN
1	C	179	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 [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	248/261 (95%)	0.24	25 (10%)	12 10	32, 49, 82, 104	0
1	B	253/261 (96%)	0.31	28 (11%)	10 8	33, 49, 86, 117	0
1	C	252/261 (96%)	-0.02	22 (8%)	16 12	25, 42, 80, 104	0
All	All	753/783 (96%)	0.18	75 (9%)	12 10	25, 47, 83, 117	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	255	ILE	8.9
1	A	164	LYS	6.4
1	A	123	VAL	6.0
1	A	232	ASP	5.4
1	B	255	ILE	5.3
1	B	192	GLU	5.2
1	A	165	ASP	5.2
1	A	193	GLU	5.0
1	C	255	ILE	4.8
1	C	192	GLU	4.8
1	B	257	ASP	4.7
1	B	256	GLU	4.6
1	A	109	GLU	4.5
1	C	256	GLU	4.3
1	C	186	SER	4.3
1	B	185	THR	4.3
1	A	124	GLU	4.2
1	B	109	GLU	4.2
1	B	258	GLU	4.2
1	B	232	ASP	4.0
1	B	108	GLN	4.0
1	B	164	LYS	4.0
1	B	259	GLU	4.0

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Mol	Chain	Res	Type	RSRZ
1	C	95	ASN	4.0
1	B	193	GLU	3.8
1	C	64	ARG	3.8
1	B	131	GLN	3.7
1	B	165	ASP	3.6
1	B	127	GLY	3.6
1	C	93	GLU	3.5
1	C	96	ALA	3.3
1	C	165	ASP	3.3
1	B	130	GLU	3.2
1	B	64	ARG	3.2
1	C	193	GLU	3.2
1	A	185	THR	3.2
1	C	191	GLU	3.2
1	A	107	ASN	3.1
1	C	164	LYS	3.1
1	C	109	GLU	3.1
1	A	125	GLN	3.1
1	A	94	ASP	2.9
1	B	125	GLN	2.9
1	A	95	ASN	2.9
1	B	107	ASN	2.9
1	C	58	ASP	2.9
1	A	44	HIS	2.8
1	A	130	GLU	2.8
1	B	97	ASP	2.7
1	B	194	ALA	2.7
1	C	59	THR	2.7
1	C	94	ASP	2.7
1	A	64	ARG	2.6
1	C	106	PRO	2.6
1	A	1	MET	2.6
1	B	243	ASP	2.5
1	C	131	GLN	2.5
1	C	107	ASN	2.5
1	C	119	MET	2.5
1	A	97	ASP	2.4
1	A	122	ASP	2.4
1	B	94	ASP	2.4
1	A	106	PRO	2.4
1	A	96	ALA	2.3
1	A	131	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	36	ASN	2.2
1	A	108	GLN	2.2
1	A	21	ASP	2.2
1	B	126	LEU	2.2
1	C	125	GLN	2.1
1	A	163	ALA	2.0
1	B	96	ALA	2.0
1	B	106	PRO	2.0
1	B	1	MET	2.0
1	C	1	MET	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.