



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:57 AM UTC

PDB ID : 6EHT / pdb_00006eht
Title : Modulation of PCNA sliding surface by p15PAF suggests a suppressive mechanism for cisplatin-induced DNA lesion bypass by pol eta holoenzyme
Authors : De March, M.; Barrera-Vilarmau, S.; Mentegari, E.; Merino, N.; Bressan, E.; Maga, G.; Crehuet, R.; Onesti, S.; Blanco, F.J.; De Biasio, A.
Deposited on : 2017-09-15
Resolution : 3.20 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:

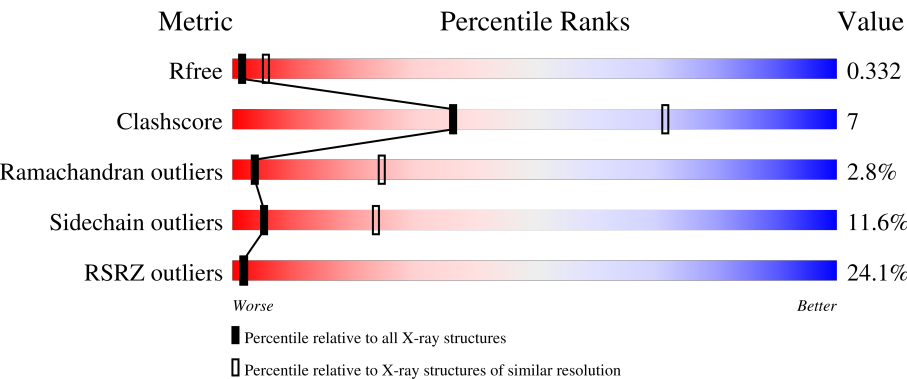
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	254	26% 77% 20% .
1	B	254	27% 83% 15% .
2	C	256	12% 69% 28% .
3	D	20	35% 80% 20%
3	E	20	60% 85% 10% 5%

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Mol	Chain	Length	Quality of chain
4	F	10	
5	G	10	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5925 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	254	Total	C	N	O	S	0	0	0
			1733	1102	289	328	14			
1	B	253	Total	C	N	O	S	0	0	0
			1652	1050	278	311	13			

- Molecule 2 is a protein called Proliferating cell nuclear antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	256	Total	C	N	O	S	0	0	0
			1844	1161	312	355	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP P12004

- Molecule 3 is a protein called PCNA-associated factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	20	Total	C	N	O	S	0	0	0
			145	98	22	24	1			
3	E	19	Total	C	N	O	S	0	0	0
			132	88	22	21	1			

- Molecule 4 is a DNA chain called DNA (5'-D(P*AP*TP*AP*CP*GP*AP*TP*GP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	10	Total	C	N	O	P	0	0	0
			210	99	42	59	10			

- Molecule 5 is a DNA chain called DNA (5'-D(P*CP*CP*CP*AP*TP*CP*GP*TP*AP*T)-

3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	G	10	Total 200	C 96	N 33	O 61	P 10	0	0	0

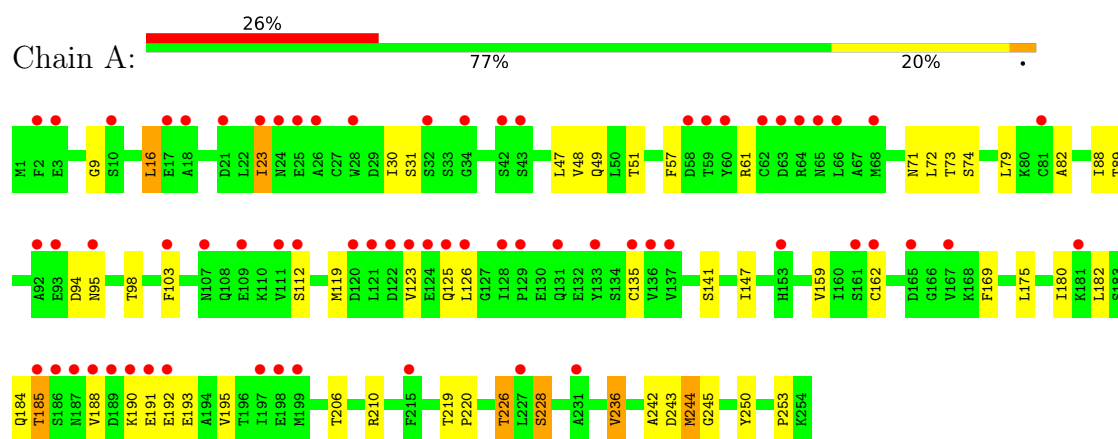
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total 3	O 3	0	0
6	B	1	Total 1	O 1	0	0
6	C	5	Total 5	O 5	0	0

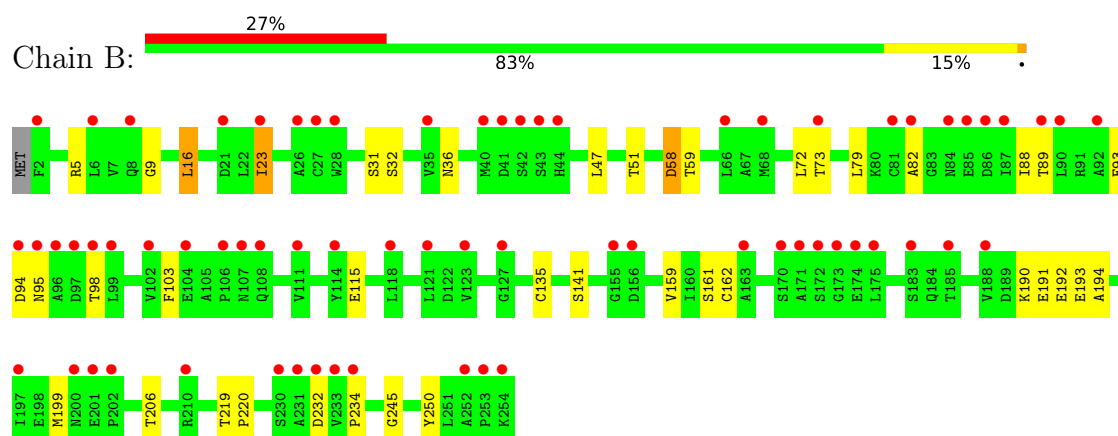
3 Residue-property plots [i](#)

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.

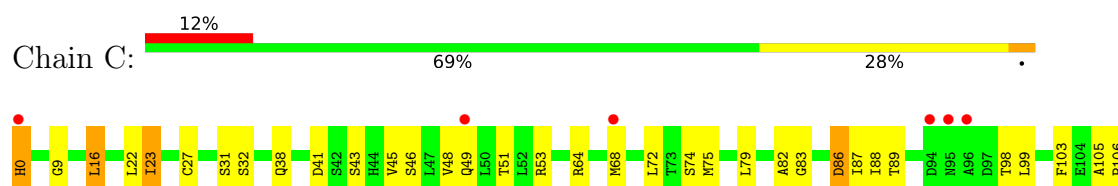
- Molecule 1: Proliferating cell nuclear antigen

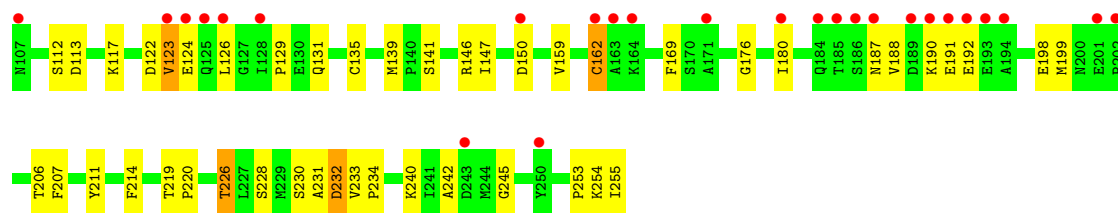


- Molecule 1: Proliferating cell nuclear antigen

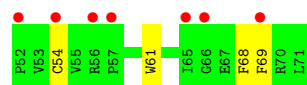
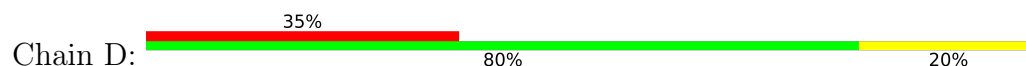


- Molecule 2: Proliferating cell nuclear antigen

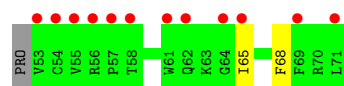
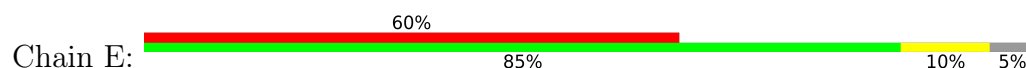




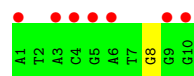
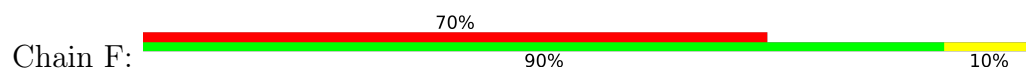
● Molecule 3: PCNA-associated factor



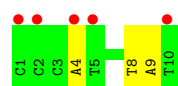
● Molecule 3: PCNA-associated factor



● Molecule 4: DNA (5'-D(P*AP*TP*AP*CP*GP*AP*TP*GP*GP*G)-3')



● Molecule 5: DNA (5'-D(P*CP*CP*CP*AP*TP*CP*GP*TP*AP*T)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.99Å 42.30Å 141.83Å 90.00° 102.70° 90.00°	Depositor
Resolution (Å)	40.48 – 3.20 40.48 – 3.20	Depositor EDS
% Data completeness (in resolution range)	94.2 (40.48-3.20) 94.2 (40.48-3.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 3.18Å)	Xtriage
Refinement program	REFMAC 5.8.0222	Depositor
R, R_{free}	0.261 , 0.327 0.259 , 0.332	Depositor DCC
R_{free} test set	717 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	70.8	Xtriage
Anisotropy	0.580	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 265.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	5925	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.22% 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	0.89	0/1759	1.07	1/2403 (0.0%)
1	B	0.88	0/1677	1.03	0/2301
2	C	1.05	2/1871 (0.1%)	1.14	0/2542
3	D	0.86	0/151	1.14	0/207
3	E	0.60	0/137	1.07	0/187
4	F	0.34	0/236	0.83	0/363
5	G	0.36	0/222	0.76	0/339
All	All	0.91	2/6053 (0.0%)	1.06	1/8342 (0.0%)

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
2	C	0	3
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	131	GLN	N-CA	-6.00	1.39	1.46
2	C	86	ASP	C-O	-5.17	1.17	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	253	PRO	CB-CA-C	5.33	117.94	110.95

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	ARG	Sidechain
2	C	0	HIS	Peptide
2	C	53	ARG	Sidechain
2	C	64	ARG	Sidechain

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	1733	0	1546	24	0
1	B	1652	0	1360	15	0
2	C	1844	0	1738	37	0
3	D	145	0	124	3	0
3	E	132	0	105	1	0
4	F	210	0	113	1	0
5	G	200	0	114	2	0
6	A	3	0	0	0	0
6	B	1	0	0	0	0
6	C	5	0	0	0	0
All	All	5925	0	5100	76	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 7.

All (76) 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:C:135:CYS:SG	2:C:162:CYS:HB3	2.26	0.76
1:A:51:THR:O	1:A:245:GLY:HA3	1.95	0.66
2:C:51:THR:O	2:C:245:GLY:HA3	1.95	0.66
5:G:8:DT:H2"	5:G:9:DA:C8	2.31	0.66
1:A:23:ILE:HB	1:A:72:LEU:HD12	1.81	0.63
2:C:45:VAL:HG11	2:C:211:TYR:CE2	2.37	0.60
1:B:58:ASP:O	1:B:59:THR:C	2.43	0.60
2:C:226:THR:O	2:C:226:THR:HG22	2.02	0.59
2:C:41:ASP:OD2	2:C:46:SER:OG	2.19	0.59
1:B:51:THR:O	1:B:245:GLY:HA3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:38:GLN:HE22	2:C:126:LEU:H	1.51	0.57
1:B:23:ILE:HB	1:B:72:LEU:HD12	1.87	0.57
2:C:23:ILE:HB	2:C:72:LEU:HD12	1.87	0.57
2:C:135:CYS:SG	2:C:162:CYS:CB	2.93	0.56
1:A:82:ALA:HB2	1:A:103:PHE:CE2	2.42	0.55
1:A:226:THR:HG22	1:A:226:THR:O	2.05	0.55
1:B:82:ALA:HB2	1:B:103:PHE:CE2	2.41	0.55
1:A:169:PHE:HE1	1:A:182:LEU:HD12	1.71	0.55
1:B:234:PRO:HD3	3:E:68:PHE:CD2	2.41	0.55
2:C:219:THR:N	2:C:220:PRO:CD	2.70	0.55
1:B:192:GLU:O	1:B:194:ALA:N	2.40	0.54
2:C:232:ASP:HB3	2:C:255:ILE:HD13	1.88	0.54
1:A:16:LEU:HD13	1:A:79:LEU:CD1	2.38	0.54
2:C:82:ALA:HB2	2:C:103:PHE:CE2	2.43	0.53
1:A:9:GLY:CA	1:A:88:ILE:HD12	2.39	0.53
1:A:219:THR:N	1:A:220:PRO:CD	2.72	0.53
1:A:184:GLN:HG3	1:A:195:VAL:O	2.09	0.53
2:C:23:ILE:HD11	2:C:48:VAL:HG11	1.91	0.53
1:A:49:GLN:NE2	1:A:51:THR:OG1	2.43	0.51
2:C:230:SER:O	2:C:231:ALA:C	2.53	0.51
2:C:16:LEU:CD1	2:C:75:MET:HE2	2.40	0.51
1:B:219:THR:N	1:B:220:PRO:CD	2.74	0.51
2:C:141:SER:HB2	2:C:219:THR:HG23	1.92	0.50
1:A:141:SER:HB2	1:A:219:THR:HG23	1.93	0.49
1:B:115:GLU:O	2:C:176:GLY:HA3	2.13	0.48
1:A:185:THR:OG1	1:A:195:VAL:N	2.42	0.48
2:C:234:PRO:HD3	3:D:68:PHE:CD2	2.49	0.48
2:C:129:PRO:HD2	3:D:69:PHE:CE1	2.49	0.48
2:C:16:LEU:HD11	2:C:75:MET:HE2	1.94	0.47
2:C:68:MET:HE1	2:C:99:LEU:HD22	1.96	0.47
4:F:8:DG:C2	5:G:4:DA:C2	3.02	0.47
2:C:232:ASP:N	2:C:232:ASP:OD1	2.49	0.46
2:C:9:GLY:CA	2:C:88:ILE:HD12	2.45	0.46
1:A:191:GLU:O	1:A:192:GLU:C	2.59	0.46
1:A:57:PHE:CE1	1:A:244:MET:HE1	2.50	0.46
2:C:135:CYS:SG	2:C:199:MET:HG3	2.56	0.46
1:B:82:ALA:HB2	1:B:103:PHE:CZ	2.51	0.45
1:A:47:LEU:HD23	1:A:250:TYR:CD1	2.52	0.45
2:C:105:ALA:HB1	2:C:106:PRO:HD2	2.00	0.44
1:B:9:GLY:CA	1:B:88:ILE:HD12	2.48	0.44
1:A:47:LEU:HB3	1:A:250:TYR:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:ILE:HG12	1:A:180:ILE:HG21	1.99	0.43
1:A:159:VAL:HG22	1:A:206:THR:OG1	2.17	0.43
1:A:71:ASN:ND2	1:A:74:SER:OG	2.52	0.43
1:A:228:SER:HB2	1:A:236:VAL:HG22	2.01	0.43
1:A:191:GLU:O	1:A:193:GLU:N	2.51	0.43
1:A:226:THR:O	1:A:226:THR:CG2	2.66	0.43
1:B:47:LEU:HD23	1:B:250:TYR:CD1	2.53	0.43
1:B:141:SER:HB2	1:B:219:THR:HG23	2.01	0.42
2:C:146:ARG:NH1	2:C:150:ASP:OD1	2.52	0.42
2:C:207:PHE:CE1	2:C:253:PRO:HB3	2.55	0.42
1:B:16:LEU:HD13	1:B:79:LEU:CD1	2.49	0.42
1:B:93:GLU:O	1:B:95:ASN:N	2.53	0.42
2:C:16:LEU:HD13	2:C:79:LEU:CD1	2.50	0.42
2:C:22:LEU:HD13	2:C:214:PHE:HB3	2.00	0.42
1:B:159:VAL:HG22	1:B:206:THR:OG1	2.19	0.42
2:C:159:VAL:HG22	2:C:206:THR:OG1	2.20	0.41
1:A:23:ILE:HD11	1:A:48:VAL:HG12	2.01	0.41
2:C:23:ILE:HD11	2:C:48:VAL:CG1	2.49	0.41
2:C:82:ALA:HB2	2:C:103:PHE:CZ	2.55	0.41
2:C:45:VAL:HG11	2:C:211:TYR:CZ	2.55	0.41
2:C:83:GLY:O	2:C:86:ASP:HB2	2.21	0.41
2:C:139:MET:HE1	2:C:169:PHE:HZ	1.84	0.41
2:C:254:LYS:HA	3:D:61:TRP:CE3	2.56	0.40
1:A:175:LEU:HD22	2:C:74:SER:HA	2.03	0.40
2:C:147:ILE:HG12	2:C:180:ILE:HG21	2.03	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	252/254 (99%)	224 (89%)	20 (8%)	8 (3%)	3 21

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	251/254 (99%)	222 (88%)	23 (9%)	6 (2%)	4	28
2	C	254/256 (99%)	232 (91%)	14 (6%)	8 (3%)	3	22
3	D	18/20 (90%)	17 (94%)	1 (6%)	0	100	100
3	E	17/20 (85%)	14 (82%)	3 (18%)	0	100	100
All	All	792/804 (98%)	709 (90%)	61 (8%)	22 (3%)	4	25

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	94	ASP
1	A	123	VAL
1	A	190	LYS
1	B	190	LYS
1	B	193	GLU
2	C	188	VAL
2	C	190	LYS
2	C	191	GLU
1	A	95	ASN
1	A	125	GLN
1	A	188	VAL
1	A	242	ALA
2	C	122	ASP
2	C	187	ASN
2	C	192	GLU
2	C	242	ALA
1	B	94	ASP
1	A	126	LEU
1	B	232	ASP
1	B	58	ASP
1	B	191	GLU
2	C	123	VAL

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	151/222 (68%)	133 (88%)	18 (12%)	5	23
1	B	124/222 (56%)	111 (90%)	13 (10%)	6	28
2	C	184/224 (82%)	161 (88%)	23 (12%)	4	21
3	D	13/18 (72%)	12 (92%)	1 (8%)	12	41
3	E	10/18 (56%)	9 (90%)	1 (10%)	7	30
All	All	482/704 (68%)	426 (88%)	56 (12%)	5	24

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	23	ILE
1	A	30	ILE
1	A	31	SER
1	A	61	ARG
1	A	73	THR
1	A	89	THR
1	A	98	THR
1	A	112	SER
1	A	119	MET
1	A	135	CYS
1	A	162	CYS
1	A	185	THR
1	A	226	THR
1	A	228	SER
1	A	236	VAL
1	A	243	ASP
1	A	244	MET
1	B	5	ARG
1	B	16	LEU
1	B	23	ILE
1	B	31	SER
1	B	32	SER
1	B	36	ASN
1	B	73	THR
1	B	89	THR
1	B	98	THR
1	B	135	CYS
1	B	161	SER
1	B	162	CYS
1	B	199	MET

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Mol	Chain	Res	Type
2	C	0	HIS
2	C	16	LEU
2	C	23	ILE
2	C	27	CYS
2	C	31	SER
2	C	32	SER
2	C	43	SER
2	C	49	GLN
2	C	87	ILE
2	C	89	THR
2	C	98	THR
2	C	112	SER
2	C	113	ASP
2	C	117	LYS
2	C	123	VAL
2	C	124	GLU
2	C	162	CYS
2	C	198	GLU
2	C	226	THR
2	C	228	SER
2	C	232	ASP
2	C	233	VAL
2	C	240	LYS
3	D	54	CYS
3	E	65	ILE

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	36	ASN
1	A	38	GLN
1	A	49	GLN
1	A	71	ASN
1	B	36	ASN
1	B	38	GLN
2	C	24	ASN
2	C	38	GLN

5.3.3 RNA ⓘ

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	254/254 (100%)	1.28	67 (26%) 1 1	30, 54, 68, 81	107 (42%)
1	B	253/254 (99%)	1.31	68 (26%) 1 1	29, 56, 70, 77	137 (54%)
2	C	256/256 (100%)	0.77	32 (12%) 8 6	28, 49, 66, 79	44 (17%)
3	D	20/20 (100%)	1.42	7 (35%) 1 1	41, 62, 88, 106	6 (30%)
3	E	19/20 (95%)	2.92	12 (63%) 0 0	29, 35, 42, 44	19 (100%)
4	F	10/10 (100%)	2.27	7 (70%) 0 0	196, 226, 243, 261	0
5	G	10/10 (100%)	2.21	5 (50%) 0 1	210, 236, 246, 270	0
All	All	822/824 (99%)	1.19	198 (24%) 2 2	28, 52, 72, 270	313 (38%)

All (198) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	55	VAL	12.3
2	C	192	GLU	9.3
2	C	191	GLU	6.1
2	C	96	ALA	5.9
1	B	233	VAL	5.8
1	B	234	PRO	5.6
2	C	126	LEU	5.6
1	A	189	ASP	5.4
1	A	191	GLU	5.4
1	B	44	HIS	5.1
2	C	202	PRO	5.1
1	B	230	SER	5.1
1	B	232	ASP	5.0
1	B	28	TRP	5.0
2	C	124	GLU	4.8
1	A	192	GLU	4.8
1	A	187	ASN	4.7

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Mol	Chain	Res	Type	RSRZ
1	A	161	SER	4.7
1	B	231	ALA	4.6
2	C	94	ASP	4.6
1	B	254	LYS	4.6
2	C	194	ALA	4.6
1	B	43	SER	4.6
3	E	65	ILE	4.6
1	B	86	ASP	4.6
2	C	193	GLU	4.5
3	D	69	PHE	4.5
1	A	186	SER	4.5
1	B	114	TYR	4.4
2	C	189	ASP	4.4
1	A	42	SER	4.2
1	A	190	LYS	4.2
2	C	184	GLN	4.1
2	C	163	ALA	4.1
1	A	198	GLU	4.1
5	G	10	DT	4.0
2	C	185	THR	4.0
1	B	107	ASN	4.0
1	B	87	ILE	4.0
1	A	185	THR	4.0
3	D	66	GLY	3.9
1	A	107	ASN	3.8
1	B	81	CYS	3.8
1	B	200	ASN	3.8
3	D	54	CYS	3.8
3	E	69	PHE	3.7
1	A	125	GLN	3.7
1	B	2	PHE	3.7
1	A	188	VAL	3.6
3	E	54	CYS	3.6
1	A	59	THR	3.6
1	B	155	GLY	3.6
1	A	43	SER	3.6
1	B	68	MET	3.5
1	A	162	CYS	3.5
1	A	199	MET	3.5
1	B	26	ALA	3.5
3	D	57	PRO	3.5
3	E	64	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	175	LEU	3.3
1	B	252	ALA	3.3
1	B	95	ASN	3.3
1	B	253	PRO	3.2
1	A	26	ALA	3.2
1	A	66	LEU	3.2
1	B	171	ALA	3.2
1	A	131	GLN	3.2
1	A	62	CYS	3.1
1	A	120	ASP	3.1
1	A	123	VAL	3.1
1	A	133	TYR	3.1
1	A	58	ASP	3.1
1	B	41	ASP	3.1
1	A	124	GLU	3.1
1	A	128	ILE	3.1
1	B	163	ALA	3.1
3	E	53	VAL	3.1
1	B	42	SER	3.0
3	E	61	TRP	3.0
4	F	6	DA	2.9
2	C	150	ASP	2.9
1	B	8	GLN	2.9
1	B	94	ASP	2.9
1	B	118	LEU	2.9
1	A	65	ASN	2.9
1	B	85	GLU	2.9
1	B	106	PRO	2.9
2	C	123	VAL	2.9
4	F	4	DC	2.9
1	A	18	ALA	2.9
1	B	92	ALA	2.9
1	B	202	PRO	2.9
1	A	111	VAL	2.9
1	A	126	LEU	2.9
1	B	73	THR	2.8
1	A	28	TRP	2.8
1	A	197	ILE	2.8
4	F	3	DA	2.8
1	B	188	VAL	2.8
1	B	84	ASN	2.8
1	B	197	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	153	HIS	2.8
1	A	32	SER	2.8
2	C	190	LYS	2.8
3	E	58	THR	2.8
1	A	215	PHE	2.8
1	B	121	LEU	2.8
1	A	10	SER	2.7
2	C	187	ASN	2.7
2	C	243	ASP	2.7
1	A	167	VAL	2.7
3	D	52	PRO	2.7
1	B	172	SER	2.7
1	B	35	VAL	2.7
2	C	107	ASN	2.7
1	A	181	LYS	2.6
1	A	103	PHE	2.6
1	A	129	PRO	2.6
1	A	63	ASP	2.6
1	B	96	ALA	2.6
1	A	3	GLU	2.6
1	B	210	ARG	2.6
1	B	23	ILE	2.5
2	C	180	ILE	2.5
3	E	57	PRO	2.5
1	B	99	LEU	2.5
1	A	24	ASN	2.5
2	C	128	ILE	2.5
1	B	174	GLU	2.5
1	B	97	ASP	2.5
5	G	5	DT	2.5
1	A	136	VAL	2.5
3	E	62	GLN	2.5
1	B	82	ALA	2.4
4	F	10	DG	2.4
1	A	121	LEU	2.4
5	G	4	DA	2.4
1	A	60	TYR	2.4
1	A	109	GLU	2.4
1	A	112	SER	2.4
2	C	162	CYS	2.4
1	B	66	LEU	2.4
2	C	49	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	183	SER	2.4
1	B	185	THR	2.4
2	C	95	ASN	2.4
1	A	21	ASP	2.4
2	C	250	TYR	2.4
5	G	1	DC	2.4
1	A	68	MET	2.4
1	B	27	CYS	2.4
1	B	21	ASP	2.4
1	B	108	GLN	2.4
1	A	2	PHE	2.4
5	G	2	DC	2.4
3	E	71	LEU	2.4
4	F	1	DA	2.4
2	C	125	GLN	2.3
1	A	122	ASP	2.3
1	B	156	ASP	2.3
1	A	25	GLU	2.3
1	B	40	MET	2.3
1	A	81	CYS	2.3
1	B	111	VAL	2.3
1	A	17	GLU	2.2
2	C	171	ALA	2.2
1	B	90	LEU	2.2
2	C	164	LYS	2.2
1	A	137	VAL	2.2
1	A	165	ASP	2.2
1	A	92	ALA	2.2
1	B	170	SER	2.2
1	B	173	GLY	2.2
1	B	98	THR	2.2
2	C	0	HIS	2.2
1	B	123	VAL	2.1
1	A	64	ARG	2.1
1	A	23	ILE	2.1
2	C	201	GLU	2.1
4	F	5	DG	2.1
1	A	93	GLU	2.1
1	A	95	ASN	2.1
1	B	102	VAL	2.1
3	D	56	ARG	2.1
1	B	104	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	6	LEU	2.1
1	B	89	THR	2.1
1	A	231	ALA	2.1
1	B	201	GLU	2.1
3	E	56	ARG	2.0
2	C	68	MET	2.0
1	A	34	GLY	2.0
1	A	227	LEU	2.0
4	F	9	DG	2.0
2	C	186	SER	2.0
3	D	65	ILE	2.0
1	A	135	CYS	2.0
1	B	127	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.