



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:36 PM GMT

PDB ID : 4CS5  
Title : Crystal Structure of PCNA from *Litopenaeus vannamei*  
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Deposited on : 2014-03-04  
Resolution : 3.00 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](mailto: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.

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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

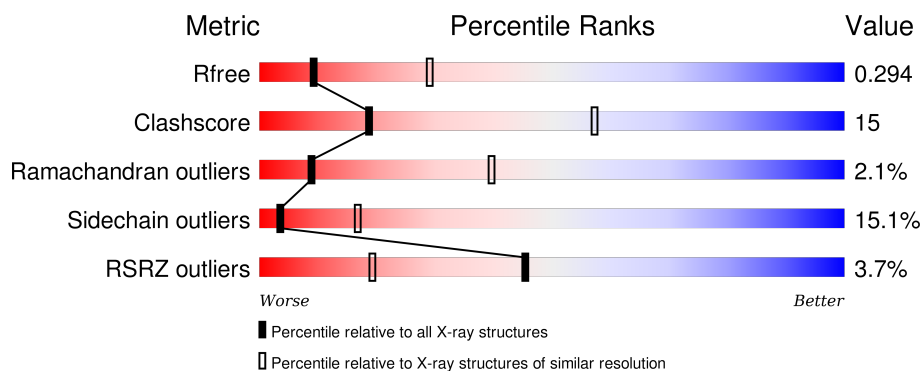
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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))
$R_{free}$	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	260	<div> <div>3%</div> <div>55%</div> <div>38%</div> <div>5%</div> </div>
1	B	260	<div> <div>4%</div> <div>56%</div> <div>36%</div> <div>5%</div> </div>
1	C	260	<div> <div>4%</div> <div>55%</div> <div>38%</div> <div>.</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5877 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.

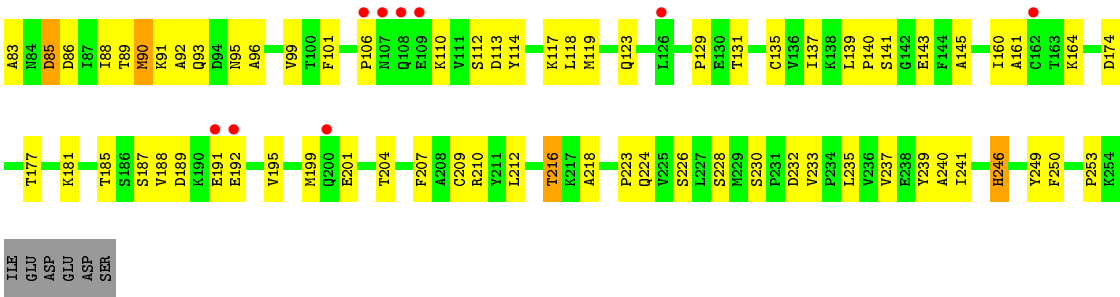
- Molecule 1 is a protein called PROLIFERATING CELL NUCLEAR ANTIGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	0	0
			1962	1239	324	380	19			
1	B	254	Total	C	N	O	S	0	0	0
			1960	1236	323	383	18			
1	C	254	Total	C	N	O	S	0	0	0
			1955	1233	321	382	19			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	55	LYS	GLU	CONFLICT	UNP G1E6N7
B	55	LYS	GLU	CONFLICT	UNP G1E6N7
C	55	LYS	GLU	CONFLICT	UNP G1E6N7





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	144.57Å 83.38Å 74.31Å 90.00° 117.65° 90.00°	Depositor
Resolution (Å)	28.06 – 3.00 28.06 – 3.00	Depositor EDS
% Data completeness (in resolution range)	93.7 (28.06-3.00) 93.7 (28.06-3.00)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.94 (at 3.00Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.265 , 0.311 0.243 , 0.294	Depositor DCC
$R_{free}$ test set	736 reflections (4.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	55.5	Xtriage
Anisotropy	0.678	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 46.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 14802 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	5877	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.51	0/1991	0.74	0/2688
1	B	0.50	0/1989	0.75	1/2685 (0.0%)
1	C	0.49	0/1983	0.72	0/2676
All	All	0.50	0/5963	0.74	1/8049 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	22	LEU	CA-CB-CG	5.67	128.34	115.30

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	1962	0	1980	65	0
1	B	1960	0	1972	60	0
1	C	1955	0	1960	59	0
All	All	5877	0	5912	182	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 15.

All (182) 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:140:PRO:HB2	1:A:143:GLU:HB2	1.70	0.73
1:B:140:PRO:HB2	1:B:143:GLU:HB2	1.70	0.73
1:C:140:PRO:HB2	1:C:143:GLU:HB2	1.72	0.72
1:A:99:VAL:HG12	1:A:118:LEU:HD11	1.72	0.72
1:C:99:VAL:HG12	1:C:118:LEU:HD11	1.72	0.71
1:A:33:SER:O	1:A:54:ALA:N	2.19	0.71
1:B:33:SER:O	1:B:54:ALA:N	2.18	0.69
1:A:21:ASP:N	1:A:21:ASP:OD1	2.23	0.68
1:B:21:ASP:N	1:B:21:ASP:OD1	2.27	0.68
1:C:145:ALA:HA	1:C:216:THR:HG21	1.77	0.67
1:B:2:PHE:HE2	1:B:90:MET:HB3	1.59	0.67
1:B:240:ALA:HA	1:B:246:HIS:HB3	1.78	0.66
1:B:145:ALA:HA	1:B:216:THR:HG21	1.77	0.65
1:B:83:ALA:N	1:B:86:ASP:OD2	2.25	0.65
1:A:145:ALA:HA	1:A:216:THR:HG21	1.79	0.65
1:A:240:ALA:HA	1:A:246:HIS:HB3	1.78	0.65
1:A:7:VAL:HG22	1:A:58:ASP:HB2	1.80	0.63
1:B:99:VAL:HG12	1:B:118:LEU:HD11	1.79	0.63
1:C:33:SER:O	1:C:54:ALA:N	2.24	0.62
1:A:2:PHE:HE2	1:A:90:MET:HB3	1.63	0.62
1:C:7:VAL:HG22	1:C:58:ASP:HB2	1.80	0.61
1:C:2:PHE:HE2	1:C:90:MET:HB3	1.65	0.61
1:C:23:LEU:HD23	1:C:41:ASP:HA	1.82	0.61
1:C:5:ARG:HG3	1:C:89:THR:OG1	2.00	0.61
1:A:241:ILE:HG13	1:A:246:HIS:HA	1.82	0.61
1:C:240:ALA:HA	1:C:246:HIS:HB3	1.82	0.61
1:B:241:ILE:HG13	1:B:246:HIS:HA	1.81	0.61
1:B:23:LEU:HD23	1:B:41:ASP:HA	1.83	0.60
1:A:23:LEU:HD23	1:A:41:ASP:HA	1.83	0.60
1:C:241:ILE:HG13	1:C:246:HIS:HA	1.82	0.60
1:C:21:ASP:N	1:C:21:ASP:OD1	2.26	0.60
1:C:83:ALA:N	1:C:86:ASP:OD2	2.29	0.59
1:C:50:LEU:HD11	1:C:52:LEU:HG	1.84	0.59
1:B:26:ALA:HB3	1:B:72:LEU:HD21	1.85	0.59
1:A:141:SER:HB2	1:A:239:TYR:HE1	1.68	0.59
1:C:141:SER:HB2	1:C:239:TYR:HE1	1.68	0.58
1:B:185:THR:OG1	1:B:195:VAL:N	2.37	0.58
1:A:85:ASP:N	1:A:85:ASP:OD1	2.37	0.58
1:B:141:SER:HB2	1:B:239:TYR:HE1	1.68	0.57
1:B:85:ASP:OD1	1:B:85:ASP:N	2.36	0.57
1:B:23:LEU:HD13	1:B:39:ALA:HB3	1.86	0.57
1:B:5:ARG:HG3	1:B:89:THR:OG1	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:LEU:HD13	1:C:39:ALA:HB3	1.86	0.57
1:B:207:PHE:HE1	1:B:253:PRO:HG3	1.70	0.56
1:C:185:THR:OG1	1:C:195:VAL:N	2.38	0.55
1:A:83:ALA:N	1:A:86:ASP:OD2	2.27	0.54
1:A:23:LEU:HD13	1:A:39:ALA:HB3	1.89	0.54
1:A:29:ASP:O	1:A:35:ILE:HA	2.07	0.54
1:A:181:LYS:O	1:C:110:LYS:HA	2.07	0.54
1:A:50:LEU:HD11	1:A:52:LEU:HG	1.89	0.54
1:B:3:GLU:HG3	1:B:91:LYS:HG2	1.88	0.54
1:A:3:GLU:HG3	1:A:91:LYS:HG2	1.90	0.54
1:C:3:GLU:HG3	1:C:91:LYS:HG2	1.88	0.53
1:C:19:ILE:O	1:C:21:ASP:N	2.42	0.53
1:A:5:ARG:HG3	1:A:89:THR:OG1	2.08	0.53
1:A:192:GLU:HG3	1:A:223:PRO:HB2	1.90	0.53
1:C:90:MET:HG2	1:C:99:VAL:HG21	1.92	0.52
1:A:207:PHE:HE1	1:A:253:PRO:HG3	1.74	0.52
1:B:37:LEU:HB3	1:B:50:LEU:HB3	1.91	0.52
1:B:192:GLU:HG3	1:B:223:PRO:HB2	1.92	0.52
1:A:2:PHE:O	1:A:91:LYS:HA	2.10	0.52
1:A:185:THR:OG1	1:A:195:VAL:N	2.39	0.52
1:B:29:ASP:O	1:B:35:ILE:HA	2.10	0.52
1:C:11:LEU:O	1:C:15:VAL:HG23	2.10	0.51
1:A:26:ALA:HB3	1:A:72:LEU:HD21	1.92	0.51
1:C:85:ASP:OD1	1:C:85:ASP:N	2.43	0.51
1:C:207:PHE:HE1	1:C:253:PRO:HG3	1.75	0.51
1:A:69:GLY:O	1:A:119:MET:HE2	2.11	0.51
1:B:137:ILE:HG22	1:B:139:LEU:HD22	1.93	0.51
1:A:137:ILE:HG22	1:A:139:LEU:HD22	1.93	0.50
1:C:29:ASP:O	1:C:35:ILE:HA	2.11	0.50
1:C:164:LYS:HD2	1:C:199:MET:HB3	1.93	0.50
1:B:164:LYS:HD2	1:B:199:MET:HB3	1.93	0.50
1:C:192:GLU:HG3	1:C:223:PRO:HB2	1.92	0.50
1:C:26:ALA:HB3	1:C:72:LEU:HD21	1.93	0.50
1:B:218:ALA:HB3	1:B:239:TYR:CD2	2.47	0.49
1:C:37:LEU:HB3	1:C:50:LEU:HB3	1.94	0.49
1:A:90:MET:HG2	1:A:99:VAL:HG21	1.95	0.49
1:B:93:GLN:H	1:B:96:ALA:HB2	1.76	0.49
1:A:12:LEU:HG	1:A:79:LEU:HD21	1.93	0.49
1:B:19:ILE:O	1:B:21:ASP:N	2.46	0.49
1:A:164:LYS:HD2	1:A:199:MET:HB3	1.95	0.48
1:C:2:PHE:O	1:C:91:LYS:HA	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:ASN:HB2	1:B:119:MET:SD	2.54	0.48
1:A:11:LEU:HA	1:A:11:LEU:HD22	1.72	0.48
1:A:19:ILE:O	1:A:21:ASP:N	2.46	0.48
1:B:34:GLY:HA2	1:B:60:TYR:CZ	2.49	0.48
1:A:218:ALA:HB3	1:A:239:TYR:CD2	2.49	0.47
1:C:218:ALA:HB3	1:C:239:TYR:CD2	2.50	0.47
1:A:10:SER:O	1:A:14:LYS:HG3	2.15	0.47
1:B:90:MET:HG2	1:B:99:VAL:HG21	1.96	0.47
1:B:93:GLN:N	1:B:96:ALA:HB2	2.30	0.47
1:A:93:GLN:H	1:A:96:ALA:HB2	1.79	0.47
1:A:11:LEU:O	1:A:15:VAL:HG23	2.15	0.47
1:C:4:ALA:HB1	1:C:57:PHE:CD2	2.50	0.47
1:C:246:HIS:ND1	1:C:246:HIS:N	2.61	0.47
1:B:9:GLY:HA2	1:B:88:ILE:HG12	1.97	0.47
1:C:93:GLN:H	1:C:96:ALA:HB2	1.80	0.46
1:B:11:LEU:O	1:B:15:VAL:HG23	2.15	0.46
1:A:53:ARG:HB2	1:A:53:ARG:HE	1.33	0.46
1:C:93:GLN:N	1:C:96:ALA:HB2	2.31	0.46
1:B:2:PHE:HA	1:B:61:ARG:O	2.16	0.46
1:B:2:PHE:O	1:B:91:LYS:HA	2.15	0.46
1:A:161:ALA:HB2	1:A:204:THR:HG23	1.98	0.46
1:B:246:HIS:ND1	1:B:246:HIS:N	2.63	0.46
1:B:160:ILE:HD12	1:B:207:PHE:CE2	2.50	0.46
1:C:160:ILE:HD12	1:C:207:PHE:CE2	2.51	0.45
1:A:71:ASN:HB2	1:A:119:MET:SD	2.56	0.45
1:A:93:GLN:N	1:A:96:ALA:HB2	2.32	0.45
1:B:4:ALA:HB1	1:B:57:PHE:CD2	2.51	0.45
1:A:112:SER:HB3	1:A:114:TYR:CE1	2.52	0.45
1:B:27:SER:HA	1:B:68:MET:O	2.17	0.45
1:B:27:SER:HB2	1:B:123:GLN:NE2	2.30	0.45
1:C:112:SER:HB3	1:C:114:TYR:CE1	2.51	0.45
1:C:137:ILE:HG22	1:C:139:LEU:HD22	1.98	0.45
1:A:2:PHE:HA	1:A:61:ARG:O	2.17	0.45
1:C:69:GLY:O	1:C:119:MET:HE2	2.16	0.45
1:B:11:LEU:HA	1:B:11:LEU:HD22	1.72	0.44
1:C:10:SER:O	1:C:14:LYS:HG3	2.17	0.44
1:A:37:LEU:HB3	1:A:50:LEU:HB3	1.99	0.44
1:A:27:SER:HA	1:A:68:MET:O	2.18	0.44
1:C:2:PHE:HA	1:C:61:ARG:O	2.18	0.44
1:A:135:CYS:SG	1:A:199:MET:HG2	2.57	0.44
1:A:16:LEU:HA	1:A:16:LEU:HD13	1.74	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:47:LEU:HB3	1:C:250:PHE:HB2	1.99	0.44
1:A:4:ALA:HB1	1:A:57:PHE:CD2	2.52	0.44
1:A:246:HIS:N	1:A:246:HIS:ND1	2.64	0.44
1:C:9:GLY:HA2	1:C:88:ILE:HG12	1.99	0.44
1:B:1:MET:HE2	1:B:1:MET:HB2	1.85	0.44
1:B:112:SER:HB3	1:B:114:TYR:CE1	2.53	0.44
1:C:135:CYS:SG	1:C:199:MET:HG2	2.58	0.43
1:B:135:CYS:SG	1:B:199:MET:HG2	2.57	0.43
1:A:241:ILE:HG22	1:A:241:ILE:O	2.18	0.43
1:A:27:SER:HB2	1:A:123:GLN:NE2	2.32	0.43
1:C:34:GLY:HA2	1:C:60:TYR:CZ	2.52	0.43
1:C:27:SER:HB2	1:C:123:GLN:NE2	2.33	0.43
1:C:1:MET:HB3	1:C:92:ALA:O	2.18	0.43
1:B:10:SER:O	1:B:14:LYS:HG3	2.19	0.43
1:A:9:GLY:HA2	1:A:88:ILE:HG12	1.99	0.43
1:B:210:ARG:O	1:B:214:MET:HE2	2.19	0.43
1:B:47:LEU:HB3	1:B:250:PHE:HB2	2.00	0.43
1:A:1:MET:HB2	1:A:1:MET:HE3	1.89	0.43
1:B:50:LEU:HD22	1:B:51:ASN:N	2.34	0.43
1:C:53:ARG:HE	1:C:53:ARG:HB2	1.34	0.43
1:A:99:VAL:CG1	1:A:118:LEU:HD11	2.46	0.42
1:A:225:VAL:HB	1:A:239:TYR:CE1	2.54	0.42
1:C:161:ALA:HB2	1:C:204:THR:HG23	2.01	0.42
1:C:4:ALA:HB1	1:C:57:PHE:CE2	2.54	0.42
1:B:69:GLY:O	1:B:119:MET:HE2	2.19	0.42
1:A:212:LEU:O	1:A:216:THR:OG1	2.38	0.42
1:A:110:LYS:HA	1:B:181:LYS:O	2.20	0.42
1:A:34:GLY:HA2	1:A:60:TYR:CZ	2.55	0.42
1:A:1:MET:HB3	1:A:92:ALA:O	2.20	0.42
1:C:191:GLU:OE1	1:C:224:GLN:NE2	2.53	0.42
1:B:19:ILE:HG23	1:B:23:LEU:HD12	2.01	0.41
1:A:57:PHE:N	1:A:57:PHE:CD1	2.88	0.41
1:B:16:LEU:HD13	1:B:16:LEU:HA	1.74	0.41
1:B:1:MET:HB3	1:B:92:ALA:O	2.20	0.41
1:B:158:ILE:HD12	1:B:212:LEU:HD12	2.01	0.41
1:C:241:ILE:O	1:C:241:ILE:HG22	2.21	0.41
1:B:50:LEU:HD21	1:B:52:LEU:HG	2.01	0.41
1:C:71:ASN:HB2	1:C:119:MET:SD	2.61	0.41
1:C:212:LEU:O	1:C:216:THR:OG1	2.38	0.41
1:C:12:LEU:O	1:C:15:VAL:HB	2.19	0.41
1:A:47:LEU:HB3	1:A:250:PHE:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:VAL:HB	1:A:249:TYR:HB2	2.02	0.41
1:A:19:ILE:HB	1:A:20:LYS:H	1.71	0.41
1:A:141:SER:HA	1:A:225:VAL:HG12	2.03	0.41
1:A:79:LEU:HD12	1:A:79:LEU:HA	1.85	0.41
1:B:57:PHE:CD1	1:B:57:PHE:N	2.89	0.41
1:B:4:ALA:HB1	1:B:57:PHE:CE2	2.55	0.41
1:A:4:ALA:HB1	1:A:57:PHE:CE2	2.56	0.41
1:B:237:VAL:HB	1:B:249:TYR:HB2	2.04	0.40
1:B:85:ASP:HB2	1:B:106:PRO:HG3	2.04	0.40
1:B:160:ILE:HD12	1:B:207:PHE:HE2	1.85	0.40
1:B:69:GLY:HA3	1:B:119:MET:O	2.21	0.40
1:C:160:ILE:HD12	1:C:207:PHE:HE2	1.85	0.40
1:C:67:ILE:O	1:C:67:ILE:HG22	2.20	0.40
1:C:237:VAL:HB	1:C:249:TYR:HB2	2.03	0.40
1:C:36:GLN:HA	1:C:50:LEU:O	2.22	0.40
1:A:160:ILE:HD12	1:A:207:PHE:CE2	2.56	0.40
1:C:85:ASP:HB2	1:C:106:PRO:HG3	2.03	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	253/260 (97%)	217 (86%)	30 (12%)	6 (2%)	7	35
1	B	252/260 (97%)	215 (85%)	32 (13%)	5 (2%)	9	41
1	C	252/260 (97%)	217 (86%)	30 (12%)	5 (2%)	9	41
All	All	757/780 (97%)	649 (86%)	92 (12%)	16 (2%)	9	40

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	LYS
1	A	129	PRO
1	A	188	VAL
1	B	20	LYS
1	B	129	PRO
1	B	188	VAL
1	C	20	LYS
1	C	129	PRO
1	C	188	VAL
1	C	54	ALA
1	A	187	SER
1	B	54	ALA
1	B	187	SER
1	C	187	SER
1	A	54	ALA
1	A	234	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	219/226 (97%)	185 (84%)	34 (16%)	3	16
1	B	219/226 (97%)	185 (84%)	34 (16%)	3	16
1	C	217/226 (96%)	186 (86%)	31 (14%)	4	19
All	All	655/678 (97%)	556 (85%)	99 (15%)	3	17

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	5	ARG
1	A	6	LEU
1	A	11	LEU
1	A	21	ASP
1	A	32	ASP
1	A	35	ILE

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Mol	Chain	Res	Type
1	A	50	LEU
1	A	79	LEU
1	A	85	ASP
1	A	90	MET
1	A	95	ASN
1	A	101	PHE
1	A	113	ASP
1	A	117	LYS
1	A	131	THR
1	A	170	SER
1	A	174	ASP
1	A	177	THR
1	A	181	LYS
1	A	189	ASP
1	A	201	GLU
1	A	209	CYS
1	A	210	ARG
1	A	216	THR
1	A	226	SER
1	A	228	SER
1	A	230	SER
1	A	232	ASP
1	A	233	VAL
1	A	235	LEU
1	A	246	HIS
1	A	254	LYS
1	A	255	ILE
1	B	1	MET
1	B	5	ARG
1	B	6	LEU
1	B	7	VAL
1	B	11	LEU
1	B	21	ASP
1	B	22	LEU
1	B	32	ASP
1	B	35	ILE
1	B	50	LEU
1	B	85	ASP
1	B	90	MET
1	B	95	ASN
1	B	101	PHE
1	B	113	ASP

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Mol	Chain	Res	Type
1	B	117	LYS
1	B	131	THR
1	B	165	GLU
1	B	170	SER
1	B	174	ASP
1	B	177	THR
1	B	181	LYS
1	B	189	ASP
1	B	209	CYS
1	B	210	ARG
1	B	216	THR
1	B	226	SER
1	B	228	SER
1	B	230	SER
1	B	232	ASP
1	B	233	VAL
1	B	235	LEU
1	B	246	HIS
1	B	254	LYS
1	C	1	MET
1	C	5	ARG
1	C	6	LEU
1	C	11	LEU
1	C	21	ASP
1	C	32	ASP
1	C	35	ILE
1	C	48	VAL
1	C	50	LEU
1	C	85	ASP
1	C	90	MET
1	C	95	ASN
1	C	101	PHE
1	C	113	ASP
1	C	117	LYS
1	C	131	THR
1	C	174	ASP
1	C	177	THR
1	C	181	LYS
1	C	189	ASP
1	C	201	GLU
1	C	209	CYS
1	C	210	ARG

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Mol	Chain	Res	Type
1	C	216	THR
1	C	226	SER
1	C	228	SER
1	C	230	SER
1	C	232	ASP
1	C	233	VAL
1	C	235	LEU
1	C	246	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	255/260 (98%)	0.07	8 (3%)	52 24	43, 67, 114, 134	0
1	B	254/260 (97%)	0.12	10 (3%)	43 18	43, 67, 113, 135	0
1	C	254/260 (97%)	0.12	10 (3%)	43 18	42, 67, 114, 134	0
All	All	763/780 (97%)	0.11	28 (3%)	45 19	42, 67, 114, 135	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	125	HIS	4.5
1	A	123	GLN	4.3
1	B	107	ASN	4.3
1	B	187	SER	4.1
1	A	120	ASN	4.0
1	A	107	ASN	3.7
1	B	106	PRO	3.3
1	C	126	LEU	3.2
1	B	186	SER	3.2
1	A	191	GLU	3.1
1	C	38	GLN	2.8
1	B	95	ASN	2.8
1	C	200	GLN	2.8
1	C	191	GLU	2.6
1	B	124	GLU	2.6
1	C	162	CYS	2.6
1	C	109	GLU	2.6
1	C	107	ASN	2.6
1	A	106	PRO	2.5
1	C	106	PRO	2.5
1	A	108	GLN	2.2
1	A	109	GLU	2.2
1	B	191	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	192	GLU	2.2
1	A	255	ILE	2.1
1	B	192	GLU	2.1
1	C	108	GLN	2.1
1	B	122	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 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.