Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 07:12 am BST

PDB ID	:	1PLR
Title	:	CRYSTAL STRUCTURE OF THE EUKARYOTIC DNA POLYMERASE PROCESSIVITY FACTOR PCNA
Authors	:	Krishna, T.S.R.; Kong, X.-P.; Gary, S.; Burgers, P.M.; Kuriyan, J.
Deposited on	:	1995-01-02
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*
A user guide is available at
https://www.wwpdb.org/validation/2017/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:

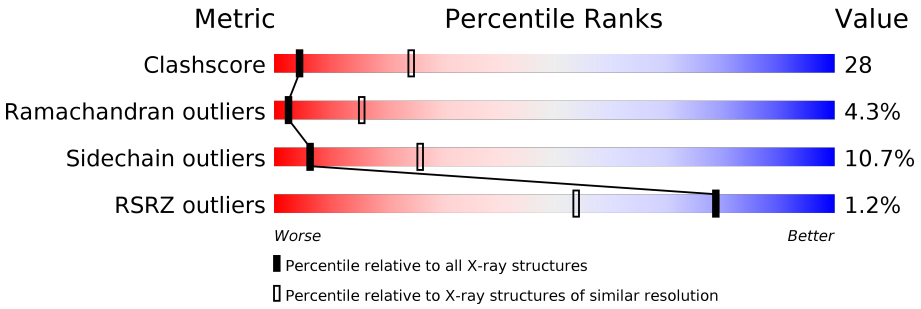
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 3.00 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))
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (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 >=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 <=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	258	% 50% 42% 7% .

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2040 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 (PCNA).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	258	2030	1295	319	406	10	0	0	0

- Molecule 2 is water.

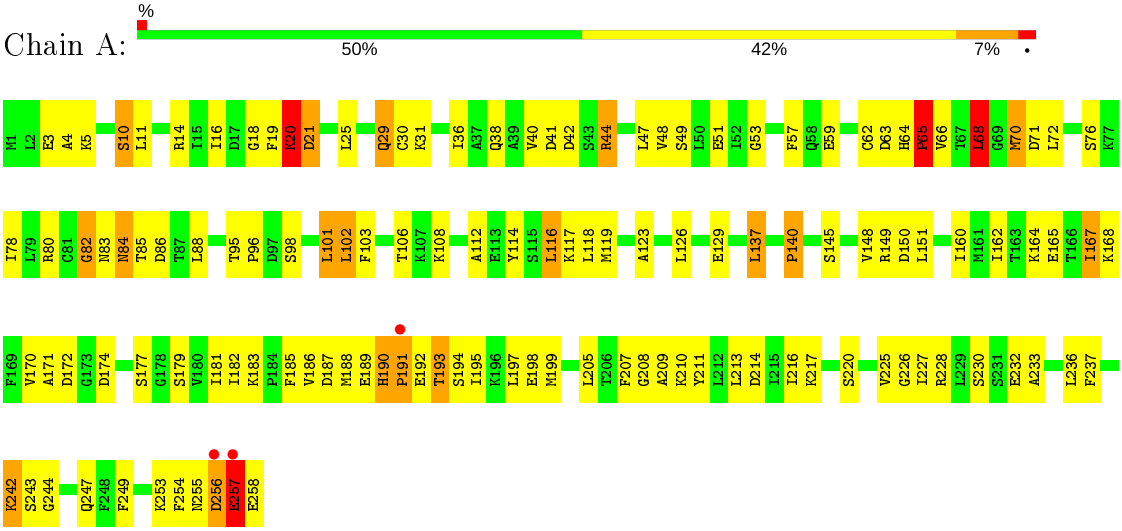
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	10	10	10	0	0

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 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 (PCNA)

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α, β, γ	121.70Å 121.70Å 121.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 3.00 38.48 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-3.00) 86.3 (38.48-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.63 (at 3.01Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.218 , (Not available) 0.213 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	56.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 77.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.062 for l,-k,h	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	2040	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.50% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.64	0/2060	0.92	6/2777 (0.2%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	68	LEU	CA-CB-CG	6.73	130.79	115.30
1	A	20	LYS	N-CA-C	-6.43	93.65	111.00
1	A	82	GLY	N-CA-C	5.94	127.94	113.10
1	A	254	PHE	N-CA-C	-5.54	96.04	111.00
1	A	126	LEU	CA-CB-CG	5.15	127.16	115.30
1	A	256	ASP	N-CA-C	-5.03	97.43	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2030	0	2041	115	0
2	A	10	0	0	0	0
All	All	2040	0	2041	115	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 28.

All (115) 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:21:ASP:HB2	1:A:217:LYS:HE3	1.42	0.98
1:A:40:VAL:HG22	1:A:47:LEU:HD12	1.43	0.98
1:A:242:LYS:HD3	1:A:242:LYS:H	1.39	0.86
1:A:98:SER:HA	1:A:118:LEU:HG	1.61	0.83
1:A:98:SER:HB3	1:A:117:LYS:HA	1.61	0.83
1:A:31:LYS:HE3	1:A:36:ILE:HG12	1.62	0.80
1:A:253:LYS:NZ	1:A:258:GLU:HG3	1.97	0.79
1:A:151:LEU:HB3	1:A:171:ALA:HB2	1.63	0.79
1:A:214:ASP:O	1:A:217:LYS:HG2	1.81	0.79
1:A:30:CYS:SG	1:A:66:VAL:HG12	2.24	0.77
1:A:253:LYS:HA	1:A:257:GLU:HB2	1.68	0.76
1:A:236:LEU:HD12	1:A:249:PHE:CE1	2.22	0.75
1:A:49:SER:HB3	1:A:247:GLN:HG3	1.69	0.75
1:A:160:ILE:HD12	1:A:207:PHE:CE1	2.24	0.72
1:A:31:LYS:NZ	1:A:31:LYS:HB2	2.06	0.70
1:A:208:GLY:HA3	1:A:258:GLU:HB3	1.74	0.68
1:A:168:LYS:HG3	1:A:181:ILE:HG12	1.76	0.67
1:A:253:LYS:HG3	1:A:256:ASP:O	1.95	0.66
1:A:188:MET:HA	1:A:191:PRO:HG3	1.77	0.66
1:A:236:LEU:HD12	1:A:249:PHE:HE1	1.60	0.66
1:A:10:SER:HB3	1:A:84:ASN:HB2	1.77	0.66
1:A:137:LEU:HD13	1:A:167:ILE:HG21	1.77	0.65
1:A:88:LEU:HD11	1:A:101:LEU:HD23	1.79	0.64
1:A:242:LYS:CD	1:A:242:LYS:H	2.03	0.63
1:A:5:LYS:O	1:A:5:LYS:HG3	2.01	0.61
1:A:25:LEU:HD22	1:A:119:MET:HG3	1.82	0.61
1:A:148:VAL:HG12	1:A:213:LEU:HD23	1.82	0.60
1:A:190:HIS:O	1:A:193:THR:HG22	2.01	0.60
1:A:20:LYS:HG3	1:A:76:SER:OG	2.02	0.59
1:A:83:ASN:O	1:A:85:THR:N	2.36	0.59
1:A:253:LYS:HZ3	1:A:258:GLU:HG3	1.69	0.58
1:A:170:VAL:HG22	1:A:179:SER:HB2	1.86	0.57
1:A:165:GLU:O	1:A:183:LYS:HD3	2.04	0.57
1:A:209:ALA:O	1:A:213:LEU:HG	2.04	0.57
1:A:14:ARG:HH21	1:A:220:SER:HB2	1.69	0.57
1:A:20:LYS:HA	1:A:72:LEU:HD13	1.85	0.57
1:A:102:LEU:O	1:A:102:LEU:HD23	2.05	0.56
1:A:214:ASP:HB3	1:A:217:LYS:NZ	2.20	0.56
1:A:129:GLU:HG3	1:A:236:LEU:CD1	2.35	0.56
1:A:256:ASP:O	1:A:257:GLU:HB2	2.07	0.55
1:A:78:ILE:HD12	1:A:116:LEU:HB2	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:LEU:O	1:A:216:ILE:HG12	2.06	0.54
1:A:145:SER:O	1:A:149:ARG:HG3	2.08	0.54
1:A:14:ARG:HH11	1:A:14:ARG:HG2	1.73	0.54
1:A:253:LYS:HD2	1:A:258:GLU:HA	1.90	0.54
1:A:208:GLY:HA3	1:A:258:GLU:CB	2.39	0.53
1:A:88:LEU:HD11	1:A:101:LEU:CD2	2.39	0.53
1:A:31:LYS:HE3	1:A:36:ILE:CG1	2.37	0.52
1:A:18:GLY:O	1:A:217:LYS:HE2	2.09	0.52
1:A:225:VAL:HG22	1:A:226:GLY:N	2.23	0.52
1:A:21:ASP:HB2	1:A:217:LYS:CE	2.26	0.52
1:A:51:GLU:O	1:A:244:GLY:HA3	2.09	0.52
1:A:38:GLN:HA	1:A:48:VAL:O	2.09	0.52
1:A:186:VAL:HG12	1:A:187:ASP:H	1.75	0.52
1:A:129:GLU:HG3	1:A:236:LEU:HD12	1.92	0.51
1:A:172:ASP:HA	1:A:177:SER:CB	2.39	0.51
1:A:29:GLN:HG2	1:A:36:ILE:HD11	1.92	0.51
1:A:253:LYS:CE	1:A:258:GLU:HG3	2.40	0.51
1:A:31:LYS:HB2	1:A:31:LYS:HZ3	1.74	0.51
1:A:228:ARG:HD2	1:A:236:LEU:HD22	1.92	0.51
1:A:232:GLU:O	1:A:233:ALA:HB2	2.11	0.51
1:A:160:ILE:HD12	1:A:207:PHE:CD1	2.46	0.50
1:A:140:PRO:HG3	1:A:193:THR:HA	1.94	0.50
1:A:208:GLY:HA2	1:A:258:GLU:HG2	1.94	0.50
1:A:4:ALA:HB1	1:A:57:PHE:CD2	2.47	0.50
1:A:151:LEU:HB3	1:A:171:ALA:CB	2.39	0.49
1:A:242:LYS:O	1:A:243:SER:HB3	2.13	0.49
1:A:25:LEU:CD2	1:A:119:MET:HG3	2.42	0.49
1:A:5:LYS:HG2	1:A:59:GLU:HB3	1.94	0.49
1:A:186:VAL:HG12	1:A:187:ASP:N	2.27	0.48
1:A:208:GLY:CA	1:A:258:GLU:HG2	2.44	0.48
1:A:98:SER:CA	1:A:118:LEU:HG	2.38	0.47
1:A:25:LEU:HD23	1:A:71:ASP:HA	1.95	0.47
1:A:20:LYS:O	1:A:21:ASP:CB	2.63	0.47
1:A:20:LYS:O	1:A:21:ASP:HB2	2.14	0.47
1:A:227:ILE:HG12	1:A:237:PHE:HD2	1.78	0.47
1:A:66:VAL:HG22	1:A:68:LEU:HD12	1.96	0.47
1:A:18:GLY:O	1:A:217:LYS:HG3	2.14	0.47
1:A:11:LEU:HD22	1:A:242:LYS:HE3	1.96	0.46
1:A:185:PHE:CE2	1:A:195:ILE:HD13	2.50	0.46
1:A:210:LYS:HG3	1:A:211:TYR:N	2.30	0.46
1:A:172:ASP:HA	1:A:177:SER:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:ILE:HG13	1:A:167:ILE:HG23	1.98	0.45
1:A:227:ILE:HG12	1:A:237:PHE:CD2	2.51	0.45
1:A:19:PHE:N	1:A:19:PHE:CD1	2.85	0.44
1:A:103:PHE:HB2	1:A:112:ALA:HB3	2.00	0.44
1:A:53:GLY:HA3	1:A:243:SER:OG	2.17	0.44
1:A:16:ILE:O	1:A:20:LYS:HB3	2.18	0.43
1:A:66:VAL:HG13	1:A:66:VAL:O	2.18	0.43
1:A:76:SER:O	1:A:80:ARG:HG3	2.19	0.43
1:A:190:HIS:O	1:A:192:GLU:N	2.52	0.43
1:A:49:SER:HB3	1:A:247:GLN:CG	2.46	0.43
1:A:205:LEU:HD23	1:A:205:LEU:HA	1.84	0.43
1:A:64:HIS:HA	1:A:65:PRO:HD2	1.94	0.43
1:A:40:VAL:HG11	1:A:44:ARG:HD3	2.00	0.42
1:A:63:ASP:O	1:A:64:HIS:CG	2.72	0.42
1:A:14:ARG:HG2	1:A:14:ARG:NH1	2.35	0.42
1:A:41:ASP:OD1	1:A:41:ASP:N	2.52	0.42
1:A:70:MET:HA	1:A:119:MET:HG2	2.01	0.42
1:A:167:ILE:HD12	1:A:182:ILE:CD1	2.50	0.42
1:A:170:VAL:HG13	1:A:179:SER:HB3	2.00	0.41
1:A:62:CYS:O	1:A:62:CYS:SG	2.77	0.41
1:A:164:LYS:HE3	1:A:165:GLU:OE1	2.21	0.41
1:A:114:TYR:CD1	1:A:114:TYR:N	2.88	0.41
1:A:170:VAL:HG22	1:A:179:SER:CB	2.49	0.41
1:A:63:ASP:O	1:A:64:HIS:ND1	2.53	0.41
1:A:189:GLU:O	1:A:190:HIS:HB2	2.20	0.41
1:A:205:LEU:HD11	1:A:230:SER:O	2.20	0.41
1:A:208:GLY:H	1:A:258:GLU:C	2.23	0.41
1:A:66:VAL:HG22	1:A:68:LEU:CD1	2.51	0.41
1:A:186:VAL:HA	1:A:194:SER:HB3	2.02	0.41
1:A:164:LYS:HA	1:A:199:MET:CE	2.51	0.41
1:A:149:ARG:HH11	1:A:149:ARG:HD2	1.75	0.40
1:A:19:PHE:CD2	1:A:48:VAL:HG11	2.57	0.40
1:A:19:PHE:CG	1:A:48:VAL:HG11	2.57	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	256/258 (99%)	212 (83%)	33 (13%)	11 (4%)	2 15

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	65	PRO
1	A	82	GLY
1	A	84	ASN
1	A	257	GLU
1	A	21	ASP
1	A	193	THR
1	A	123	ALA
1	A	150	ASP
1	A	190	HIS
1	A	255	ASN
1	A	191	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	233/233 (100%)	208 (89%)	25 (11%)	6 26

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLU
1	A	10	SER
1	A	20	LYS
1	A	29	GLN
1	A	42	ASP
1	A	44	ARG
1	A	65	PRO
1	A	68	LEU
1	A	70	MET
1	A	86	ASP
1	A	95	THR
1	A	96	PRO
1	A	101	LEU
1	A	102	LEU
1	A	106	THR
1	A	108	LYS
1	A	116	LEU
1	A	137	LEU
1	A	140	PRO
1	A	167	ILE
1	A	174	ASP
1	A	197	LEU
1	A	198	GLU
1	A	242	LYS
1	A	257	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	GLN
1	A	247	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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	258/258 (100%)	-0.43	3 (1%) 79 54	4, 29, 77, 89	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	256	ASP	3.0
1	A	257	GLU	2.6
1	A	191	PRO	2.5

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.