



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2024 – 11:20 AM EDT

PDB ID : 6K00  
Title : Crystal structure A of ceNAP1-H2A-H2B complex  
Authors : Liu, Y.R.  
Deposited on : 2019-05-05  
Resolution : 2.20 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](mailto: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>.

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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
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

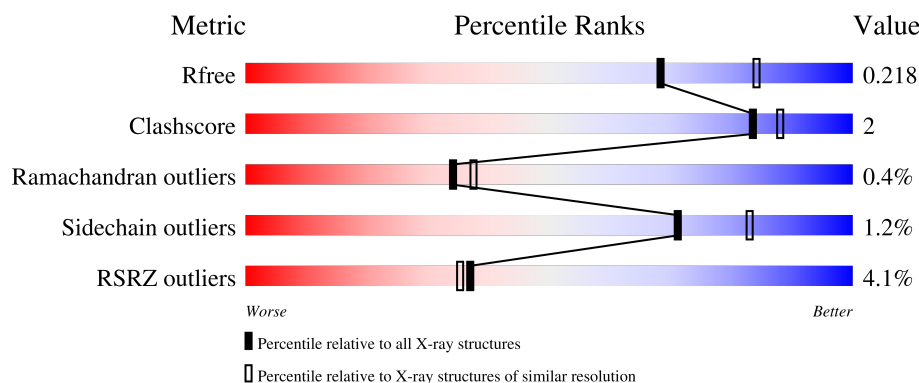
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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  $\geq 3$ , 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	308	
1	B	308	
2	D	207	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5636 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 Nucleosome Assembly Protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2083	1341	336	402	4			
1	B	269	Total	C	N	O	S	0	0	0
			2136	1370	346	415	5			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	initiating methionine	UNP Q19007
A	-10	GLY	-	expression tag	UNP Q19007
A	-9	SER	-	expression tag	UNP Q19007
A	-8	SER	-	expression tag	UNP Q19007
A	-7	HIS	-	expression tag	UNP Q19007
A	-6	HIS	-	expression tag	UNP Q19007
A	-5	HIS	-	expression tag	UNP Q19007
A	-4	HIS	-	expression tag	UNP Q19007
A	-3	HIS	-	expression tag	UNP Q19007
A	-2	HIS	-	expression tag	UNP Q19007
A	-1	SER	-	expression tag	UNP Q19007
A	0	SER	-	expression tag	UNP Q19007
A	1	GLY	-	expression tag	UNP Q19007
A	2	GLU	-	expression tag	UNP Q19007
A	3	ASN	-	expression tag	UNP Q19007
A	4	LEU	-	expression tag	UNP Q19007
A	5	TYR	-	expression tag	UNP Q19007
A	6	PHE	-	expression tag	UNP Q19007
A	7	GLN	-	expression tag	UNP Q19007
A	8	HIS	-	expression tag	UNP Q19007
A	9	MET	-	expression tag	UNP Q19007
B	-11	MET	-	initiating methionine	UNP Q19007
B	-10	GLY	-	expression tag	UNP Q19007
B	-9	SER	-	expression tag	UNP Q19007
B	-8	SER	-	expression tag	UNP Q19007

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP Q19007
B	-6	HIS	-	expression tag	UNP Q19007
B	-5	HIS	-	expression tag	UNP Q19007
B	-4	HIS	-	expression tag	UNP Q19007
B	-3	HIS	-	expression tag	UNP Q19007
B	-2	HIS	-	expression tag	UNP Q19007
B	-1	SER	-	expression tag	UNP Q19007
B	0	SER	-	expression tag	UNP Q19007
B	1	GLY	-	expression tag	UNP Q19007
B	2	GLU	-	expression tag	UNP Q19007
B	3	ASN	-	expression tag	UNP Q19007
B	4	LEU	-	expression tag	UNP Q19007
B	5	TYR	-	expression tag	UNP Q19007
B	6	PHE	-	expression tag	UNP Q19007
B	7	GLN	-	expression tag	UNP Q19007
B	8	HIS	-	expression tag	UNP Q19007
B	9	MET	-	expression tag	UNP Q19007

- Molecule 2 is a protein called Histone H2B 1,Histone H2A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	180	Total	C	N	O	S	0	0	0
			1369	858	255	254	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	28	HIS	-	expression tag	UNP P04255
D	29	MET	-	expression tag	UNP P04255

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	25	Total	O	0	0
			25	25		
3	B	19	Total	O	0	0
			19	19		
3	D	4	Total	O	0	0
			4	4		



- Molecule 1: Nucleosome Assembly Protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	68.30Å 78.72Å 169.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.35 – 2.20 49.35 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.7 (49.35-2.20) 98.7 (49.35-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.76 (at 2.20Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, $R_{free}$	0.211 , 0.243 0.213 , 0.218	Depositor DCC
$R_{free}$ test set	2365 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.8	Xtriage
Anisotropy	0.421	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 37.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5636	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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, 0.333 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	0.32	0/2127	0.52	2/2881 (0.1%)
1	B	0.28	0/2182	0.46	0/2954
2	D	0.25	0/1388	0.41	0/1875
All	All	0.29	0/5697	0.47	2/7710 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	152	ASP	CB-CG-OD1	-6.93	112.06	118.30
1	A	190	GLN	CB-CA-C	-5.30	99.80	110.40

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	2083	0	2044	11	1
1	B	2136	0	2066	10	1
2	D	1369	0	1388	8	0
3	A	25	0	0	0	0
3	B	19	0	0	0	0
3	D	4	0	0	0	0
All	All	5636	0	5498	27	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 2.

All (27) 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:187:ALA:O	1:A:190:GLN:HG3	1.76	0.86
1:B:253:ARG:HH12	1:B:263:GLU:HG2	1.47	0.79
1:B:250:LYS:HE2	1:B:253:ARG:HH21	1.52	0.74
1:A:264:GLU:O	1:A:268:LEU:HD12	1.90	0.70
1:A:247:PRO:HG2	1:A:265:PHE:HB2	1.79	0.64
1:B:248:LYS:HD2	1:B:248:LYS:H	1.68	0.56
1:A:261:GLN:H	1:A:261:GLN:CD	2.09	0.56
1:A:63:LYS:HE3	1:A:67:ASP:OD2	2.06	0.55
1:B:215:LYS:HG3	1:B:235:VAL:HG23	1.89	0.55
2:D:86:ILE:HG13	2:D:155:ALA:HB3	1.90	0.54
1:A:262:ALA:HA	1:A:265:PHE:CD1	2.44	0.52
1:A:196:VAL:HG21	1:A:270:TYR:CG	2.49	0.48
1:B:145:VAL:HG12	1:B:281:ILE:HD11	1.97	0.47
2:D:75:SER:HB2	2:D:154:TYR:HB3	1.98	0.46
2:D:84:SER:OG	2:D:84:SER:O	2.30	0.45
1:B:250:LYS:O	1:B:250:LYS:HD3	2.16	0.45
2:D:87:SER:OG	2:D:88:SER:N	2.49	0.45
2:D:67:PHE:CD2	2:D:166:LEU:HD21	2.53	0.44
2:D:189:LYS:HB2	2:D:189:LYS:HE2	1.69	0.44
1:A:283:ARG:NH2	1:B:191:PHE:O	2.51	0.43
1:A:95:ASP:O	1:A:99:GLU:HG2	2.18	0.43
2:D:109:SER:O	2:D:113:LYS:HG3	2.18	0.43
1:B:292:LEU:HD23	1:B:292:LEU:HA	1.87	0.43
1:B:175:LYS:HE3	1:B:246:PRO:HB3	2.02	0.42
1:A:96:GLN:NE2	1:B:11:LEU:HD11	2.34	0.41
2:D:195:PRO:O	2:D:199:GLN:HG2	2.21	0.41
1:A:264:GLU:O	1:A:268:LEU:CD1	2.63	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:A:63:LYS:NZ	1:B:67:ASP:OD2[3_544]	2.15	0.05

## 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	257/308 (83%)	254 (99%)	3 (1%)	0	100	100
1	B	263/308 (85%)	260 (99%)	3 (1%)	0	100	100
2	D	176/207 (85%)	169 (96%)	4 (2%)	3 (2%)	9	6
All	All	696/823 (85%)	683 (98%)	10 (1%)	3 (0%)	34	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	101	GLY
2	D	157	ARG
2	D	154	TYR

### 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	A	222/265 (84%)	220 (99%)	2 (1%)	78	88
1	B	226/265 (85%)	223 (99%)	3 (1%)	69	81
2	D	140/166 (84%)	138 (99%)	2 (1%)	67	80
All	All	588/696 (84%)	581 (99%)	7 (1%)	71	83

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	LYS
1	A	218	LYS
1	B	45	SER
1	B	248	LYS
1	B	294	SER
2	D	75	SER
2	D	203	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	156	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 monosaccharides 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	263/308 (85%)	0.27	10 (3%) 40 38	23, 36, 70, 132	0
1	B	269/308 (87%)	0.10	11 (4%) 37 35	26, 37, 78, 111	0
2	D	180/207 (86%)	0.13	8 (4%) 34 32	29, 44, 96, 131	0
All	All	712/823 (86%)	0.17	29 (4%) 37 35	23, 39, 82, 132	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	262	ALA	9.6
1	A	260	GLU	4.5
2	D	155	ALA	3.8
1	A	265	PHE	3.7
2	D	156	GLN	3.5
1	A	231	LEU	3.4
1	A	261	GLN	3.1
1	B	265	PHE	3.0
1	A	190	GLN	3.0
2	D	203	ARG	2.9
1	A	296	ASP	2.8
1	B	91	GLY	2.8
1	B	217	VAL	2.8
1	A	263	GLU	2.6
1	B	252	GLU	2.6
2	D	77	LEU	2.6
2	D	78	ALA	2.5
1	B	92	LEU	2.5
1	B	110	ALA	2.5
2	D	85	THR	2.4
1	A	233	LYS	2.4
1	A	268	LEU	2.4
1	B	251	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	186	GLU	2.3
1	B	261	GLN	2.3
1	B	108	PRO	2.1
2	D	122	LYS	2.0
1	B	185	ALA	2.0
2	D	129	ALA	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 monosaccharides 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.