

wwPDB X-ray Structure Validation Summary Report (i)

Nov 20, 2023 – 12:17 AM JST

PDB ID	:	7C7X
Title	:	Structural insights into nucleosome reorganization by NAP1-RELATED PRO-
		TEIN 1 (NRP1)
Authors	:	Luo, Q.; Baihui, W.
Deposited on	:	2020-05-27
Resolution	:	3.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (1)) were used in the production of this report:

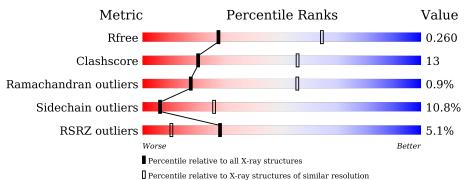
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\;DIFFRACTION$

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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R _{free}	130704	2092 (3.00-3.00)
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 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 <=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 chai	in			
1	А	93	% 62%	2:	3%	•	13%
1	С	93	<u>4%</u> 57%	28%)	·	13%
2	В	98	74%		14%	•	7%
2	D	98	69%		17%	5%	8%
3	Е	239		13% •		27%	
3	F	239	8% 49% 17%	•	3	31%	



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5502 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 Histone H2A.6.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace	
1	С	81	Total 620		N 114	O 110	0	0	0
1	А	81	Total 615	-	N 112	O 110	0	0	0

• Molecule 2 is a protein called Histone H2B.1.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
0	Л	90	Total	С	Ν	0	S	0	0	0
	D	90	707	455	120	130	2	0	0	0
0	р	91	Total	С	Ν	0	S	0	0	0
	D	91	715	459	122	132	2		0	0

• Molecule 3 is a protein called NAP1-related protein 1.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	Е	174	Total	С	Ν	0	S	0	0	0
5	Ľ	174	1426	922	221	282	1	0	0	0
2	Б	166	Total	С	Ν	0	S	0	0	0
5	Г	100	1371	890	211	269	1	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

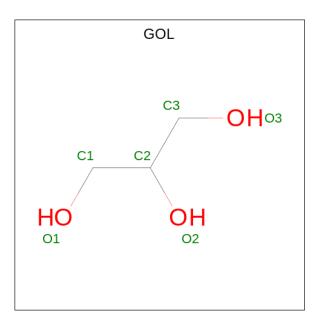
Chain	Residue	Modelled	Actual	Comment	Reference
Е	18	SER	-	expression tag	UNP Q9CA59
F	18	SER	-	expression tag	UNP Q9CA59

• Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).









Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
4	D	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0
4	D	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0
4	F	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0

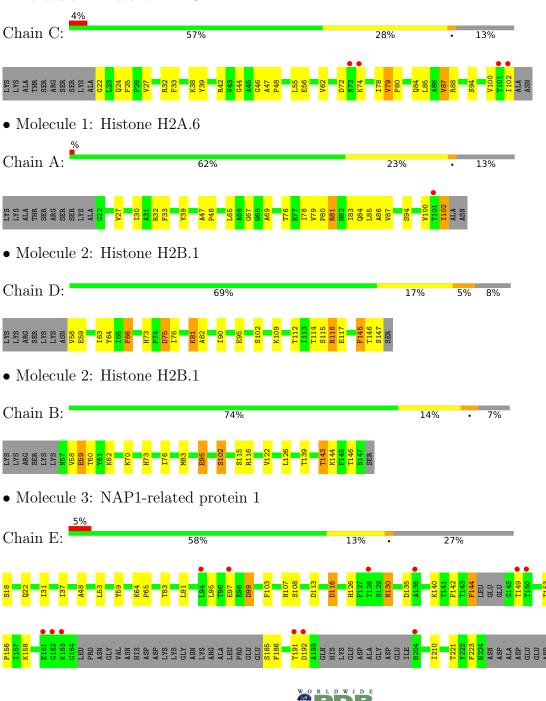
• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	С	4	Total O 4 4	0	0
5	D	3	Total O 3 3	0	0
5	Ε	3	Total O 3 3	0	0
5	F	5	$\begin{array}{cc} \text{Total} & \text{O} \\ 5 & 5 \end{array}$	0	0
5	А	6	Total O 6 6	0	0
5	В	3	Total O 3 3	0	0



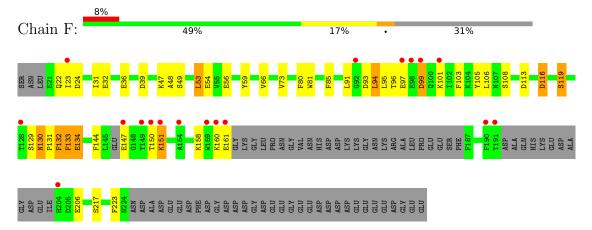
3 Residue-property plots (i)

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: Histone H2A.6

• Molecule 3: NAP1-related protein 1





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	66.71Å 128.40Å 140.28Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.62 - 3.00	Depositor
Resolution (A)	29.60 - 2.99	EDS
% Data completeness	84.9 (29.62-3.00)	Depositor
(in resolution range)	84.2 (29.60-2.99)	EDS
R _{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$3.10 (at 3.00 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
D D.	0.223 , 0.262	Depositor
R, R_{free}	0.227 , 0.260	DCC
R_{free} test set	1044 reflections $(4.94%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	54.9	Xtriage
Anisotropy	0.201	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.32, 68.1	EDS
L-test for twinning ²	$ \langle L \rangle = 0.39, \langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5502	wwPDB-VP
Average B, all atoms $(Å^2)$	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.97% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	Bo	nd lengths	Bond angles		
MOI	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.73	1/622~(0.2%)	0.85	0/840	
1	С	0.70	0/628	0.84	0/848	
2	В	0.70	0/726	0.87	0/976	
2	D	0.71	0/718	0.90	0/965	
3	Е	0.71	0/1457	0.86	1/1969~(0.1%)	
3	F	0.67	0/1401	0.82	0/1895	
All	All	0.70	1/5552~(0.0%)	0.85	1/7493~(0.0%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
1	А	69	ALA	C-O	5.64	1.34	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	Е	192	ASP	CB-CG-OD2	5.16	122.94	118.30

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.



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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	615	0	647	31	0
1	С	620	0	652	21	0
2	В	715	0	753	11	0
2	D	707	0	747	23	0
3	Ε	1426	0	1382	19	0
3	F	1371	0	1333	55	0
4	D	18	0	24	2	0
4	F	6	0	8	0	0
5	А	6	0	0	2	0
5	В	3	0	0	0	0
5	С	4	0	0	1	0
5	D	3	0	0	1	0
5	Е	3	0	0	1	0
5	F	5	0	0	2	0
All	All	5502	0	5546	139	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 13.

The worst 5 of 139 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)
3:F:132:PHE:O	3:F:160:LYS:N	1.81	1.14
3:F:131:PRO:O	3:F:160:LYS:HB3	1.51	1.10
3:F:131:PRO:O	3:F:160:LYS:CB	1.99	1.09
1:A:81:ARG:HD2	1:A:102:ILE:HD11	1.33	1.09
3:F:95:LEU:HD11	3:F:103:PHE:CE2	1.93	1.02

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	Favoured Allowed		Percentiles
1	А	79/93~(85%)	69~(87%)	10 (13%)	0	100 100
1	С	79/93~(85%)	65~(82%)	11 (14%)	3~(4%)	3 18
2	В	89/98~(91%)	83~(93%)	5~(6%)	1 (1%)	14 50
2	D	88/98~(90%)	81 (92%)	7 (8%)	0	100 100
3	Ε	166/239~(70%)	141 (85%)	23~(14%)	2(1%)	13 48
3	F	158/239~(66%)	133 (84%)	25 (16%)	0	100 100
All	All	659/860~(77%)	572 (87%)	81 (12%)	6 (1%)	17 55

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	Ε	223	PHE
1	С	100	VAL
1	С	74	LYS
2	В	70	LYS
3	Е	37	ILE

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
1	А	60/70~(86%)	56~(93%)	4 (7%)	16 49		
1	С	61/70~(87%)	56~(92%)	5 (8%)	11 39		
2	В	80/87~(92%)	71~(89%)	9 (11%)	6 24		
2	D	79/87~(91%)	70~(89%)	9 (11%)	5 24		
3	Ε	159/214~(74%)	144 (91%)	15~(9%)	8 32		
3	F	153/214~(72%)	131~(86%)	22~(14%)	3 15		
All	All	592/742~(80%)	528~(89%)	64 (11%)	6 26		

 $5~{\rm of}~64$ residues with a non-rotameric side chain are listed below:

Mol	Chain	\mathbf{Res}	Type
2	В	60	THR

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Mol	Chain	Res	Type
2	В	102	SER
3	Е	130	ASN
3	Е	116	ASP
2	В	115	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
3	Ε	130	ASN
3	F	100	GLN
3	F	130	ASN
3	Е	22	GLN
2	D	73	HIS

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)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).



Mol	Iol Type Chain Res		s Link	Bond lengths			B	Bond ang	gles		
	туре	Chain	nes	nes	nes Lilik	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
4	GOL	F	301	-	5,5,5	0.57	0	$5,\!5,\!5$	0.47	0	
4	GOL	D	201	-	$5,\!5,\!5$	0.09	0	$5,\!5,\!5$	0.37	0	
4	GOL	D	202	-	$5,\!5,\!5$	0.58	0	$5,\!5,\!5$	0.36	0	
4	GOL	D	203	-	$5,\!5,\!5$	0.65	0	$5,\!5,\!5$	0.45	0	

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	F	301	-	-	4/4/4/4	-
4	GOL	D	201	-	-	2/4/4/4	-
4	GOL	D	202	-	-	0/4/4/4	-
4	GOL	D	203	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	201	GOL	O1-C1-C2-C3
4	F	301	GOL	C1-C2-C3-O3
4	F	301	GOL	O1-C1-C2-C3
4	D	201	GOL	O1-C1-C2-O2
4	F	301	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Μ	ol	Chain	Res	Type	Clashes	Symm-Clashes
4	=	D	201	GOL	2	0

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	$\langle RSRZ \rangle$	$\# RSRZ {>}2$	$\mathbf{OWAB}(\mathbf{\AA}^2)$	Q < 0.9
1	А	81/93~(87%)	-0.27	1 (1%) 79 54	25,65,118,144	0
1	С	81/93~(87%)	-0.22	4 (4%) 29 11	25, 63, 119, 154	0
2	В	91/98~(92%)	-0.53	0 100 100	25, 52, 94, 105	0
2	D	90/98~(91%)	-0.54	0 100 100	26, 52, 93, 115	0
3	Ε	174/239~(72%)	0.13	12 (6%) 16 5	39, 89, 156, 184	0
3	F	166/239~(69%)	0.24	18 (10%) 5 2	43, 93, 167, 190	0
All	All	683/860~(79%)	-0.11	35 (5%) 28 10	25, 71, 152, 190	0

The worst 5 of 35 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	147	GLU	6.2
3	Е	161	GLU	4.8
3	Е	162	GLY	4.2
1	С	73	ASN	4.1
3	F	150	THR	3.9

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)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
4	GOL	D	203	6/6	0.83	0.21	$28,\!31,\!31,\!31$	0
4	GOL	D	202	6/6	0.92	0.27	28,30,30,30	0
4	GOL	D	201	6/6	0.94	0.22	$53,\!57,\!59,\!61$	0
4	GOL	F	301	6/6	0.95	0.30	30,32,32,32	0

6.5 Other polymers (i)

There are no such residues in this entry.

