



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 22, 2018 – 06:28 PM EST

PDB ID : 6EQT
Title : CRYSTAL STRUCTURE OF THE HUMAN KINETOCHORE PROTEIN CENP-N
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Deposited on : 2017-10-15
Resolution : 2.73 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

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20030736

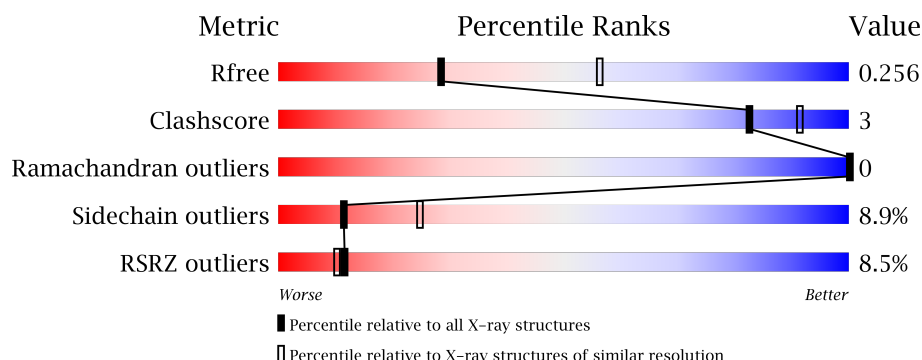
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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}	100719	3342 (2.78-2.70)
Clashscore	112137	3731 (2.78-2.70)
Ramachandran outliers	110173	3670 (2.78-2.70)
Sidechain outliers	110143	3671 (2.78-2.70)
RSRZ outliers	101464	3362 (2.78-2.70)

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. 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	219	<div> <div>7%</div> <div> <div></div> <div>79%</div> <div>15%</div> <div>• 5%</div> </div> </div>
1	B	219	<div> <div>9%</div> <div> <div></div> <div>79%</div> <div>14%</div> <div>• 6%</div> </div> </div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 3478 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 Centromere protein N.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	1	0
			1741	1122	301	310	8			
1	B	205	Total	C	N	O	S	0	0	0
			1716	1105	298	305	8			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	84	ASP	GLU	variant	UNP Q96H22
A	214	HIS	-	expression tag	UNP Q96H22
A	215	HIS	-	expression tag	UNP Q96H22
A	216	HIS	-	expression tag	UNP Q96H22
A	217	HIS	-	expression tag	UNP Q96H22
A	218	HIS	-	expression tag	UNP Q96H22
A	219	HIS	-	expression tag	UNP Q96H22
B	84	ASP	GLU	variant	UNP Q96H22
B	214	HIS	-	expression tag	UNP Q96H22
B	215	HIS	-	expression tag	UNP Q96H22
B	216	HIS	-	expression tag	UNP Q96H22
B	217	HIS	-	expression tag	UNP Q96H22
B	218	HIS	-	expression tag	UNP Q96H22
B	219	HIS	-	expression tag	UNP Q96H22

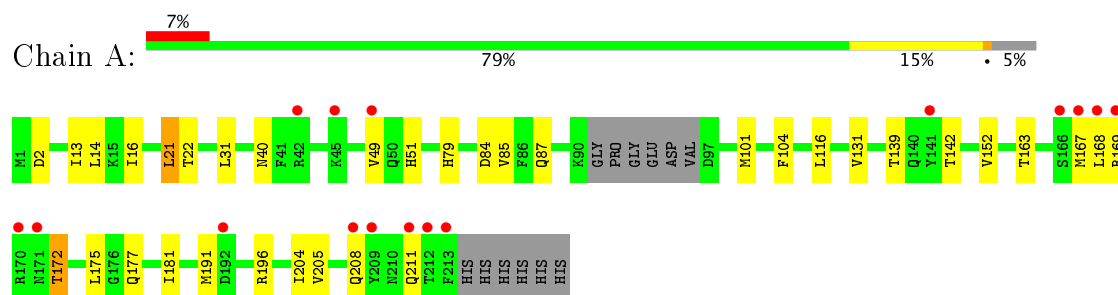
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	13	Total	O	0	0
			13	13		
2	B	8	Total	O	0	0
			8	8		

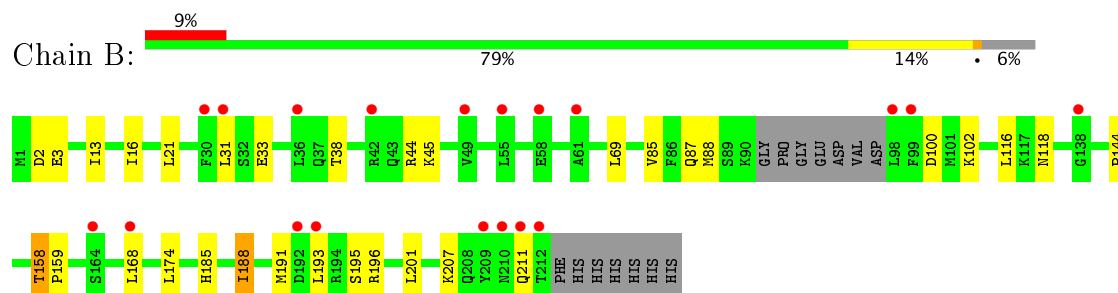
3 Residue-property plots [i](#)

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: Centromere protein N



• Molecule 1: Centromere protein N



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	87.30Å 87.30Å 81.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	87.30 – 2.73 49.12 – 2.74	Depositor EDS
% Data completeness (in resolution range)	99.8 (87.30-2.73) 99.8 (49.12-2.74)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.206 , 0.254 0.207 , 0.256	Depositor DCC
R_{free} test set	812 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	82.1	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 60.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.030 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3478	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.58% 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	0.42	0/1782	0.62	0/2403
1	B	0.42	0/1753	0.64	0/2364
All	All	0.42	0/3535	0.63	0/4767

There are no bond length outliers.

There are no bond angle outliers.

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	1741	0	1767	12	0
1	B	1716	0	1746	8	0
2	A	13	0	0	0	0
2	B	8	0	0	0	0
All	All	3478	0	3513	20	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 3.

All (20) 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:B:88:MET:HG3	1:B:188:ILE:HD12	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:THR:HG22	1:B:159:PRO:HD2	1.79	0.63
1:A:177:GLN:O	1:A:181:ILE:HG12	1.99	0.61
1:B:118:ASN:HB3	1:B:144:PRO:HD3	1.88	0.56
1:B:87:GLN:HB2	1:B:191:MET:HE2	1.87	0.55
1:A:168:LEU:HG	1:A:172:THR:HG21	1.89	0.55
1:A:87:GLN:HB2	1:A:191:MET:HE2	1.93	0.51
1:A:152:VAL:HG22	1:A:163:THR:HG22	1.94	0.49
1:A:85:VAL:HG11	1:A:205:VAL:HG11	1.94	0.49
1:A:13:ILE:HG23	1:A:21:LEU:HD21	1.94	0.48
1:A:79:HIS:O	1:A:196:ARG:HG2	2.13	0.48
1:A:208:GLN:HA	1:A:211:GLN:HG2	1.96	0.47
1:B:85:VAL:HG22	1:B:191:MET:HE1	1.97	0.47
1:A:13:ILE:HA	1:A:16:ILE:HD12	1.99	0.45
1:B:88:MET:HB3	1:B:185:HIS:CD2	2.53	0.43
1:B:195:SER:HB3	1:B:201:LEU:HG	2.00	0.43
1:B:13:ILE:HA	1:B:16:ILE:HD12	2.00	0.42
1:A:101:MET:HE3	1:A:104:PHE:HB3	2.01	0.42
1:A:14:LEU:HD23	1:A:49:VAL:HG11	2.01	0.42
1:A:139:THR:OG1	1:A:142:THR:HG22	2.20	0.41

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	204/219 (93%)	198 (97%)	6 (3%)	0	100	100
1	B	201/219 (92%)	193 (96%)	8 (4%)	0	100	100
All	All	405/438 (92%)	391 (96%)	14 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	194/203 (96%)	180 (93%)	14 (7%)	17	36
1	B	191/203 (94%)	171 (90%)	20 (10%)	8	17
All	All	385/406 (95%)	351 (91%)	34 (9%)	11	25

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	21	LEU
1	A	22	THR
1	A	31	LEU
1	A	40	ASN
1	A	51	HIS
1	A	84	ASP
1	A	116	LEU
1	A	131	VAL
1	A	167	MET
1	A	169	ARG
1	A	172	THR
1	A	175	LEU
1	A	204	ILE
1	B	2	ASP
1	B	3	GLU
1	B	21	LEU
1	B	31	LEU
1	B	33	GLU
1	B	38	THR
1	B	44	ARG
1	B	45	LYS
1	B	69	LEU
1	B	100	ASP
1	B	102	LYS
1	B	116	LEU
1	B	158	THR

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Mol	Chain	Res	Type
1	B	168	LEU
1	B	174	LEU
1	B	188	ILE
1	B	193	LEU
1	B	196	ARG
1	B	207	LYS
1	B	211	GLN

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	19	ASN
1	B	51	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 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, 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	207/219 (94%)	0.59	16 (7%) 14 13	55, 81, 134, 175	0
1	B	205/219 (93%)	0.58	19 (9%) 9 8	60, 88, 125, 171	0
All	All	412/438 (94%)	0.58	35 (8%) 11 10	55, 85, 128, 175	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	211	GLN	5.8
1	A	168	LEU	5.7
1	B	209	TYR	5.2
1	B	212	THR	4.8
1	A	167	MET	4.7
1	B	98	LEU	4.2
1	A	141	TYR	4.1
1	A	170	ARG	3.9
1	B	99	PHE	3.8
1	A	45	LYS	3.4
1	B	30	PHE	3.2
1	A	211	GLN	3.2
1	B	31	LEU	3.2
1	A	169	ARG	3.2
1	A	171	ASN	3.1
1	B	58	GLU	3.0
1	B	210	ASN	2.9
1	A	209	TYR	2.9
1	A	166	SER	2.9
1	B	138	GLY	2.8
1	A	49	VAL	2.8
1	A	42	ARG	2.8
1	A	192	ASP	2.5
1	A	213	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	36	LEU	2.5
1	B	193	LEU	2.5
1	B	192	ASP	2.4
1	B	164	SER	2.3
1	B	168	LEU	2.3
1	B	42	ARG	2.2
1	B	55	LEU	2.2
1	A	212	THR	2.1
1	B	49	VAL	2.0
1	B	61	ALA	2.0
1	A	208	GLN	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.