



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 05:17 pm BST

PDB ID : 4PXT
Title : Structural basis for the assembly of the mitotic motor kinesin-5 into bipolar tetramers
Authors : Scholey, J.E.; Nithianantham, S.; Scholey, J.M.; Al-Bassam, J.
Deposited on : 2014-03-25
Resolution : 2.90 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

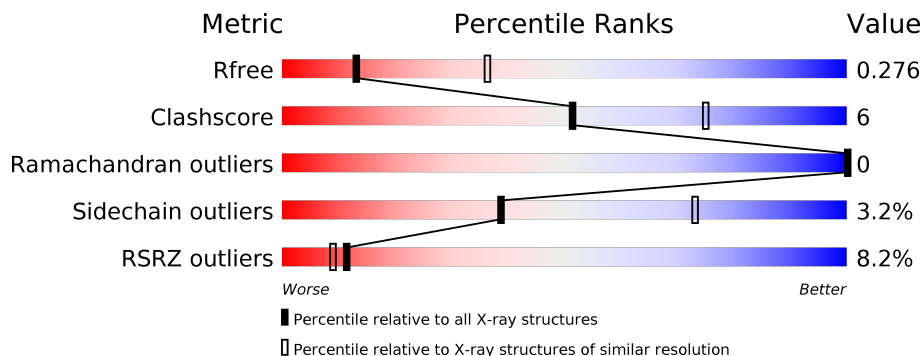
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 $\leq 5\%$. 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	211	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 65%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 11%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 23%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">5% 65% 11% • 23%</p>
1	B	211	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 63%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 27%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">7% 63% 10% • 27%</p>

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 2518 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 Bipolar kinesin KRP-130.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	163	1291	781	229	267	1	13	0	0	0
1	B	155	1227	743	218	253	1	12	0	0	0

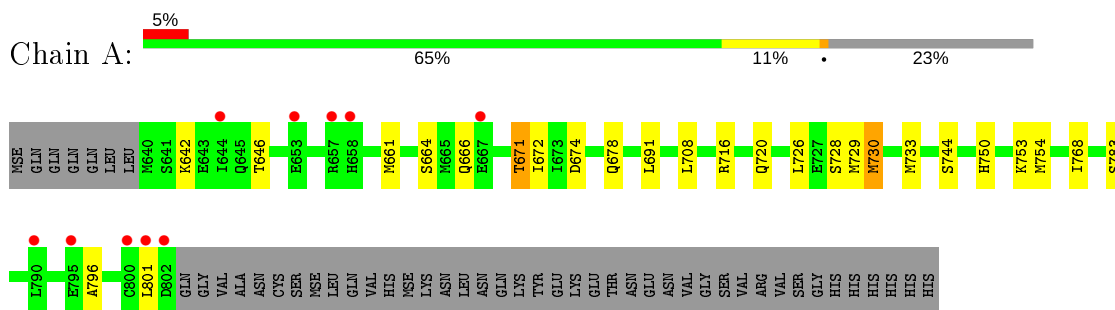
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	633	MSE	-	EXPRESSION TAG	UNP P46863
A	836	SER	HIS	ENGINEERED MUTATION	UNP P46863
A	837	GLY	HIS	ENGINEERED MUTATION	UNP P46863
A	840	HIS	-	EXPRESSION TAG	UNP P46863
A	841	HIS	-	EXPRESSION TAG	UNP P46863
A	840	HIS	-	EXPRESSION TAG	UNP P46863
A	841	HIS	-	EXPRESSION TAG	UNP P46863
A	842	HIS	-	EXPRESSION TAG	UNP P46863
A	843	HIS	-	EXPRESSION TAG	UNP P46863
B	633	MSE	-	EXPRESSION TAG	UNP P46863
B	836	SER	HIS	ENGINEERED MUTATION	UNP P46863
B	837	GLY	HIS	ENGINEERED MUTATION	UNP P46863
B	840	HIS	-	EXPRESSION TAG	UNP P46863
B	841	HIS	-	EXPRESSION TAG	UNP P46863
B	840	HIS	-	EXPRESSION TAG	UNP P46863
B	841	HIS	-	EXPRESSION TAG	UNP P46863
B	842	HIS	-	EXPRESSION TAG	UNP P46863
B	843	HIS	-	EXPRESSION TAG	UNP P46863

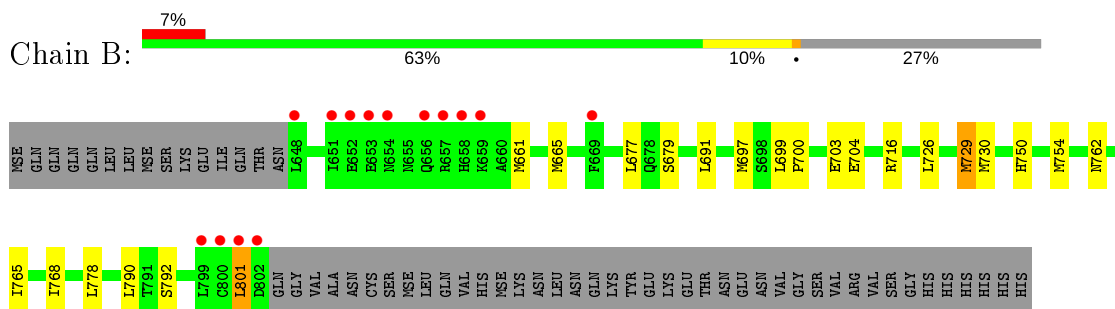
3 Residue-property plots [i](#)

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: Bipolar kinesin KRP-130



- Molecule 1: Bipolar kinesin KRP-130



4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	139.07Å 139.07Å 104.35Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	60.22 – 2.90 69.54 – 2.90	Depositor EDS
% Data completeness (in resolution range)	92.4 (60.22-2.90) 92.4 (69.54-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.80 (at 2.91Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.242 , 0.274 0.245 , 0.276	Depositor DCC
R_{free} test set	610 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	57.6	Xtrriage
Anisotropy	0.047	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2518	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.28% 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 [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	0.29	0/1286	0.40	0/1696
1	B	0.32	0/1222	0.54	1/1611 (0.1%)
All	All	0.31	0/2508	0.47	1/3307 (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	801	LEU	CA-CB-CG	5.81	128.67	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	1291	0	1283	19	0
1	B	1227	0	1218	19	0
All	All	2518	0	2501	28	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 6.

All (28) 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:691:LEU:HD22	1:B:765:ILE:HD11	1.80	0.64
1:A:733:MSE:HE2	1:A:733:MSE:HA	1.82	0.62
1:B:750:HIS:O	1:B:754:MSE:HG3	2.00	0.62
1:A:671:THR:HG22	1:A:672:ILE:HD12	1.86	0.56
1:A:691:LEU:HD11	1:B:768:ILE:HG22	1.87	0.56
1:A:729:MSE:HE1	1:B:729:MSE:HE1	1.87	0.55
1:A:768:ILE:HG22	1:B:691:LEU:HD11	1.88	0.55
1:A:666:GLN:HG3	1:B:790:LEU:HD11	1.89	0.54
1:A:796:ALA:HA	1:A:801:LEU:HD13	1.90	0.54
1:A:750:HIS:O	1:A:754:MSE:HG3	2.10	0.51
1:B:661:MSE:O	1:B:665:MSE:HE3	2.13	0.48
1:A:666:GLN:CG	1:B:790:LEU:HD11	2.44	0.48
1:A:783:SER:OG	1:B:677:LEU:HB2	2.14	0.48
1:A:730:MSE:HE1	1:B:726:LEU:HG	1.96	0.48
1:B:699:LEU:O	1:B:703:GLU:HG2	2.15	0.47
1:A:726:LEU:HD13	1:B:730:MSE:HE1	1.97	0.47
1:A:744:SER:OG	1:B:716:ARG:HD3	2.15	0.46
1:A:716:ARG:O	1:A:720:GLN:HG3	2.16	0.45
1:B:765:ILE:HA	1:B:765:ILE:HD12	1.83	0.45
1:B:762:ASN:O	1:B:765:ILE:HG22	2.16	0.45
1:A:661:MSE:HA	1:A:664:SER:HB3	1.99	0.44
1:B:726:LEU:HA	1:B:726:LEU:HD12	1.79	0.43
1:B:697:MSE:C	1:B:700:PRO:HD2	2.39	0.42
1:B:762:ASN:HA	1:B:765:ILE:HG22	2.02	0.42
1:A:674:ASP:O	1:A:678:GLN:HG2	2.20	0.41
1:A:642:LYS:O	1:A:646:THR:HG23	2.21	0.41
1:B:700:PRO:O	1:B:704:GLU:HG2	2.20	0.41
1:A:753:LYS:HB2	1:A:753:LYS:HE2	1.88	0.40

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	161/211 (76%)	158 (98%)	3 (2%)	0	100	100
1	B	153/211 (72%)	151 (99%)	2 (1%)	0	100	100
All	All	314/422 (74%)	309 (98%)	5 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	146/174 (84%)	142 (97%)	4 (3%)	44	77
1	B	138/174 (79%)	133 (96%)	5 (4%)	35	69
All	All	284/348 (82%)	275 (97%)	9 (3%)	39	73

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	671	THR
1	A	708	LEU
1	A	728	SER
1	A	730	MSE
1	B	679	SER
1	B	729	MSE
1	B	778	LEU
1	B	792	SER
1	B	801	LEU

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

Mol	Chain	Res	Type
1	B	666	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 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	150/211 (71%)	0.27	10 (6%) 17 13	33, 80, 152, 175	0
1	B	143/211 (67%)	0.31	14 (9%) 7 5	39, 75, 149, 204	0
All	All	293/422 (69%)	0.29	24 (8%) 11 9	33, 77, 152, 204	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	800	CYS	5.2
1	B	656	GLN	3.7
1	A	801	LEU	3.7
1	A	802	ASP	3.7
1	B	801	LEU	3.2
1	A	658	HIS	3.1
1	B	659	LYS	3.1
1	B	799	LEU	3.1
1	B	648	LEU	3.1
1	B	653	GLU	3.0
1	B	654	ASN	2.7
1	A	790	LEU	2.5
1	B	658	HIS	2.5
1	B	802	ASP	2.5
1	A	653	GLU	2.4
1	A	795	GLU	2.3
1	A	644	ILE	2.3
1	B	669	PHE	2.2
1	A	657	ARG	2.2
1	B	651	ILE	2.1
1	B	657	ARG	2.1
1	A	667	GLU	2.1
1	A	800	CYS	2.1
1	B	652	GLU	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.