Full wwPDB X-ray Structure Validation Report (i)

May 14, 2020 – 01:34 pm BST

PDB ID	:	3CEQ
Title	:	The TPR domain of Human Kinesin Light Chain 2 (hKLC2)
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Deposited on	:	2008-02-29
Resolution	:	2.75 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

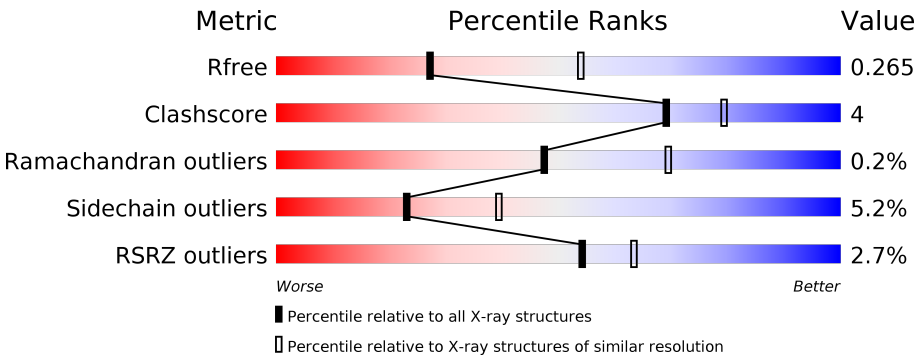
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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	283	2% 76% 11% • 12%
1	B	283	3% 81% 13% • 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4087 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 Kinesin light chain 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	250	1961	1223	358	372	8	0	0	0
			Total	C	N	O	S			
1	B	269	2124	1325	386	405	8	0	0	0

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	198	MET	-	EXPRESSION TAG	UNP Q9H0B6
A	199	GLY	-	EXPRESSION TAG	UNP Q9H0B6
A	200	SER	-	EXPRESSION TAG	UNP Q9H0B6
A	201	SER	-	EXPRESSION TAG	UNP Q9H0B6
A	202	HIS	-	EXPRESSION TAG	UNP Q9H0B6
A	203	HIS	-	EXPRESSION TAG	UNP Q9H0B6
A	204	HIS	-	EXPRESSION TAG	UNP Q9H0B6
A	205	HIS	-	EXPRESSION TAG	UNP Q9H0B6
A	206	HIS	-	EXPRESSION TAG	UNP Q9H0B6
A	207	HIS	-	EXPRESSION TAG	UNP Q9H0B6
A	208	SER	-	EXPRESSION TAG	UNP Q9H0B6
A	209	SER	-	EXPRESSION TAG	UNP Q9H0B6
A	210	GLY	-	EXPRESSION TAG	UNP Q9H0B6
A	211	LEU	-	EXPRESSION TAG	UNP Q9H0B6
A	212	VAL	-	EXPRESSION TAG	UNP Q9H0B6
A	213	PRO	-	EXPRESSION TAG	UNP Q9H0B6
A	214	ARG	-	EXPRESSION TAG	UNP Q9H0B6
A	215	GLY	-	EXPRESSION TAG	UNP Q9H0B6
A	216	SER	-	EXPRESSION TAG	UNP Q9H0B6
B	198	MET	-	EXPRESSION TAG	UNP Q9H0B6
B	199	GLY	-	EXPRESSION TAG	UNP Q9H0B6
B	200	SER	-	EXPRESSION TAG	UNP Q9H0B6
B	201	SER	-	EXPRESSION TAG	UNP Q9H0B6
B	202	HIS	-	EXPRESSION TAG	UNP Q9H0B6
B	203	HIS	-	EXPRESSION TAG	UNP Q9H0B6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	204	HIS	-	EXPRESSION TAG	UNP Q9H0B6
B	205	HIS	-	EXPRESSION TAG	UNP Q9H0B6
B	206	HIS	-	EXPRESSION TAG	UNP Q9H0B6
B	207	HIS	-	EXPRESSION TAG	UNP Q9H0B6
B	208	SER	-	EXPRESSION TAG	UNP Q9H0B6
B	209	SER	-	EXPRESSION TAG	UNP Q9H0B6
B	210	GLY	-	EXPRESSION TAG	UNP Q9H0B6
B	211	LEU	-	EXPRESSION TAG	UNP Q9H0B6
B	212	VAL	-	EXPRESSION TAG	UNP Q9H0B6
B	213	PRO	-	EXPRESSION TAG	UNP Q9H0B6
B	214	ARG	-	EXPRESSION TAG	UNP Q9H0B6
B	215	GLY	-	EXPRESSION TAG	UNP Q9H0B6
B	216	SER	-	EXPRESSION TAG	UNP Q9H0B6

- Molecule 2 is water.

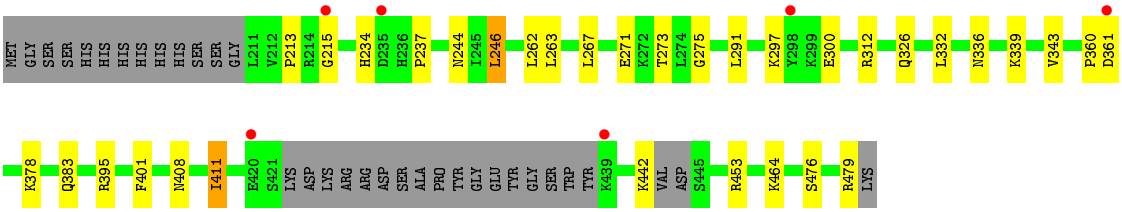
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	B	2	2	2	0	0

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.

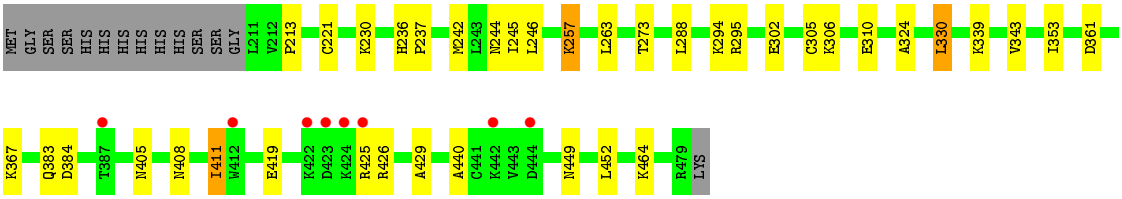
- Molecule 1: Kinesin light chain 2

Chain A:



- Molecule 1: Kinesin light chain 2

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	70.42Å 99.94Å 103.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.75 29.51 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.4 (30.00-2.75) 98.4 (29.51-2.75)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.84 (at 2.76Å)	Xtriage
Refinement program	REFMAC refmac_5.2.0019	Depositor
R, R_{free}	0.234 , 0.271 0.230 , 0.265	Depositor DCC
R_{free} test set	979 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å^2)	85.8	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 70.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.004 for -h,l,k	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4087	wwPDB-VP
Average B, all atoms (Å^2)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.04% 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 (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	0.31	0/1991	0.47	1/2682 (0.0%)
1	B	0.31	0/2162	0.48	1/2916 (0.0%)
All	All	0.31	0/4153	0.47	2/5598 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	213	PRO	N-CA-CB	5.89	110.37	103.30
1	A	213	PRO	N-CA-CB	5.88	110.36	103.30

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	1961	0	1956	13	0
1	B	2124	0	2107	17	0
2	B	2	0	0	0	0
All	All	4087	0	4063	30	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 4.

All (30) 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:405:ASN:H	1:B:408:ASN:HB3	1.49	0.76
1:A:360:PRO:HB2	1:A:395:ARG:HD3	1.81	0.62
1:A:246:LEU:HB3	1:A:262:LEU:HD13	1.83	0.60
1:A:297:LYS:HE2	1:A:300:GLU:HG3	1.85	0.59
1:B:237:PRO:HD3	1:B:273:THR:HG21	1.86	0.56
1:B:221:CYS:HB2	1:B:246:LEU:HD13	1.91	0.53
1:B:425:ARG:HD3	1:B:429:ALA:HB3	1.89	0.53
1:A:263:LEU:HD12	1:A:291:LEU:HD12	1.91	0.53
1:B:302:GLU:HG2	1:B:306:LYS:HE2	1.91	0.52
1:B:257:LYS:HD3	1:B:295:ARG:NE	2.26	0.51
1:B:305:CYS:SG	1:B:330:LEU:HD22	2.51	0.50
1:A:453:ARG:HH22	1:A:476:SER:HB3	1.77	0.49
1:B:419:GLU:HG3	1:B:464:LYS:HB3	1.94	0.49
1:A:411:ILE:H	1:A:411:ILE:HD12	1.78	0.49
1:A:361:ASP:HB3	1:A:395:ARG:HH21	1.80	0.47
1:B:263:LEU:HB3	1:B:288:LEU:HD13	1.98	0.46
1:B:236:HIS:CG	1:B:237:PRO:HD2	2.50	0.46
1:A:237:PRO:HB3	1:A:273:THR:HG21	1.97	0.45
1:B:242:MET:HA	1:B:245:ILE:HD12	1.98	0.45
1:B:449:ASN:HA	1:B:452:LEU:HD12	2.00	0.44
1:B:324:ALA:HB2	1:B:353:ILE:HB	1.99	0.43
1:B:257:LYS:HD3	1:B:295:ARG:CZ	2.50	0.42
1:A:332:LEU:O	1:A:336:ASN:ND2	2.47	0.42
1:A:312:ARG:NH1	1:A:326:GLN:HG2	2.35	0.42
1:A:401:PHE:O	1:A:408:ASN:ND2	2.53	0.42
1:B:411:ILE:HD12	1:B:440:ALA:HB1	2.02	0.42
1:A:271:GLU:O	1:A:275:GLY:HA2	2.21	0.41
1:B:339:LYS:O	1:B:343:VAL:HG23	2.21	0.41
1:A:339:LYS:O	1:A:343:VAL:HG23	2.21	0.41
1:B:221:CYS:CB	1:B:246:LEU:HD13	2.50	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	244/283 (86%)	234 (96%)	9 (4%)	1 (0%)	34	53
1	B	267/283 (94%)	260 (97%)	7 (3%)	0	100	100
All	All	511/566 (90%)	494 (97%)	16 (3%)	1 (0%)	47	69

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	215	GLY

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	203/236 (86%)	193 (95%)	10 (5%)	25	43
1	B	220/236 (93%)	208 (94%)	12 (6%)	21	37
All	All	423/472 (90%)	401 (95%)	22 (5%)	23	39

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	234	HIS
1	A	244	ASN
1	A	246	LEU
1	A	267	LEU
1	A	378	LYS
1	A	383	GLN
1	A	411	ILE
1	A	442	LYS
1	A	464	LYS
1	A	479	ARG
1	B	230	LYS
1	B	244	ASN
1	B	257	LYS
1	B	294	LYS
1	B	310	GLU

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Mol	Chain	Res	Type
1	B	330	LEU
1	B	361	ASP
1	B	367	LYS
1	B	383	GLN
1	B	384	ASP
1	B	411	ILE
1	B	426	ARG

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

Mol	Chain	Res	Type
1	A	408	ASN

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 (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, 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	250/283 (88%)	0.34	6 (2%)	59	68	71, 77, 83, 85	0
1	B	269/283 (95%)	0.30	8 (2%)	50	59	67, 77, 80, 84	0
All	All	519/566 (91%)	0.32	14 (2%)	54	63	67, 77, 82, 85	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	424	LYS	4.2
1	B	442	LYS	3.9
1	B	425	ARG	3.6
1	A	298	TYR	3.1
1	B	444	ASP	2.8
1	A	235	ASP	2.6
1	A	420	GLU	2.5
1	B	387	THR	2.5
1	A	215	GLY	2.5
1	A	361	ASP	2.5
1	B	423	ASP	2.4
1	B	412	TRP	2.3
1	A	439	LYS	2.3
1	B	422	LYS	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.