



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:21 AM UTC

PDB ID : 5W9A / pdb_00005w9a
Title : The structure of the Trim5alpha Bbox- coiled coil in complex LC3B
Authors : Keown, J.R.; Goldstone, D.C.
Deposited on : 2017-06-22
Resolution : 2.74 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

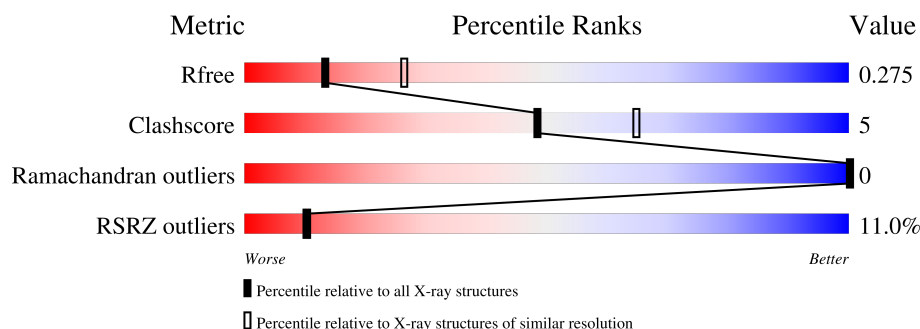
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 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}	180053	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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 $\leq 5\%$. 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	193	
1	B	193	
2	C	114	
2	D	114	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4849 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 Tripartite motif-containing protein 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1522	952	258	301	11			
1	B	191	Total	C	N	O	S	0	0	0
			1504	941	251	300	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	120	LYS	GLU	engineered mutation	UNP Q0PF16
A	121	ASP	ARG	engineered mutation	UNP Q0PF16
B	120	LYS	GLU	engineered mutation	UNP Q0PF16
B	121	ASP	ARG	engineered mutation	UNP Q0PF16

- Molecule 2 is a protein called Microtubule-associated proteins 1A/1B light chain 3B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	114	Total	C	N	O	S	0	0	0
			911	583	160	166	2			
2	D	113	Total	C	N	O	S	0	0	0
			908	580	159	167	2			

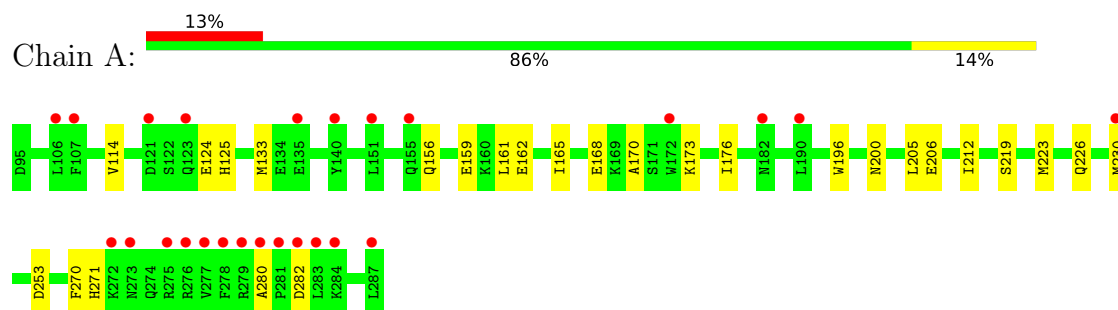
- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	2	Total	Zn	0	0
			2	2		

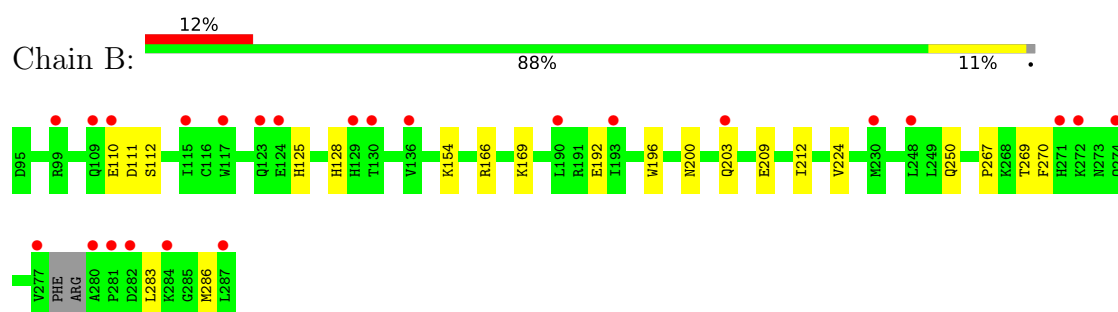
3 Residue-property plots

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.

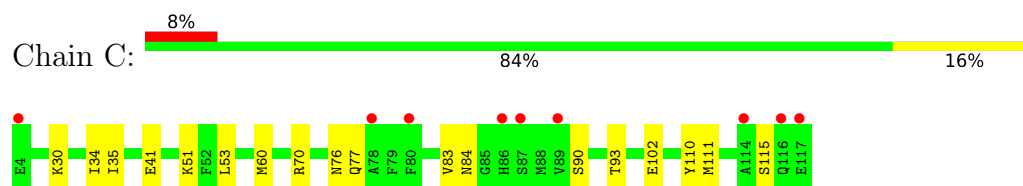
- Molecule 1: Tripartite motif-containing protein 5



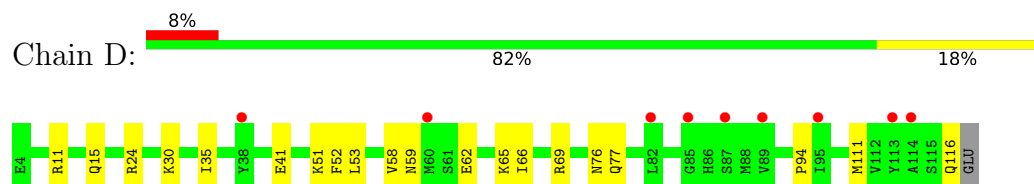
- Molecule 1: Tripartite motif-containing protein 5



- Molecule 2: Microtubule-associated proteins 1A/1B light chain 3B



- Molecule 2: Microtubule-associated proteins 1A/1B light chain 3B



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.01Å 115.32Å 174.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.91 – 2.74 40.91 – 2.74	Depositor EDS
% Data completeness (in resolution range)	62.3 (40.91-2.74) 62.3 (40.91-2.74)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.98 (at 2.73Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.260 , 0.274 0.258 , 0.275	Depositor DCC
R_{free} test set	1214 reflections (3.12%)	wwPDB-VP
Wilson B-factor (Å ²)	83.7	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 44.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4849	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.95% 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](#)

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

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	0.13	0/1543	0.28	0/2077
1	B	0.13	0/1525	0.25	0/2054
2	C	0.09	0/928	0.29	0/1256
2	D	0.09	0/923	0.29	0/1246
All	All	0.12	0/4919	0.28	0/6633

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 [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	1522	0	1445	23	0
1	B	1504	0	1414	19	0
2	C	911	0	898	13	0
2	D	908	0	912	15	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
All	All	4849	0	4669	52	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 5.

All (52) 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)
2:D:30:LYS:NZ	1:B:200:ASN:OD1	2.32	0.61
2:D:66:ILE:HG12	2:D:69:ARG:HH12	1.65	0.61
2:D:41:GLU:OE1	2:D:116:GLN:NE2	2.33	0.60
2:C:41:GLU:OE2	2:C:115:SER:N	2.35	0.59
2:C:84:ASN:ND2	2:C:102:GLU:O	2.32	0.56
2:D:76:ASN:OD1	2:D:77:GLN:N	2.39	0.55
2:D:11:ARG:HD2	1:B:192:GLU:OE2	2.07	0.55
1:A:161:LEU:HD21	1:B:267:PRO:HD3	1.89	0.55
2:D:35:ILE:HG12	2:D:111:MET:HB2	1.88	0.53
2:C:60:MET:HE1	2:C:111:MET:HE1	1.90	0.53
1:A:253:ASP:OD2	1:B:112:SER:HB2	2.08	0.53
2:C:76:ASN:OD1	2:C:77:GLN:N	2.43	0.52
2:D:53:LEU:HD12	1:B:200:ASN:HB2	1.92	0.52
2:C:83:VAL:HG22	2:C:111:MET:HG2	1.93	0.51
1:B:283:LEU:HA	1:B:286:MET:HE3	1.93	0.51
1:A:170:ALA:O	1:A:173:LYS:HG2	2.11	0.51
2:D:59:ASN:HA	2:D:94:PRO:HA	1.93	0.50
1:A:200:ASN:HB2	2:C:53:LEU:HD12	1.94	0.50
1:A:200:ASN:OD1	2:C:30:LYS:NZ	2.44	0.49
1:A:165:ILE:HD12	1:B:267:PRO:HB3	1.94	0.49
1:A:230:MET:HE2	1:B:154:LYS:HB3	1.95	0.48
1:A:162:GLU:OE2	1:B:224:VAL:HG12	2.13	0.48
1:A:212:ILE:HG23	1:A:270:PHE:HE1	1.79	0.48
1:A:270:PHE:O	1:A:271:HIS:ND1	2.39	0.48
1:A:168:GLU:OE2	1:B:269:THR:OG1	2.26	0.47
2:D:53:LEU:O	1:B:203:GLN:NE2	2.46	0.47
2:C:34:ILE:HB	2:C:110:TYR:HD1	1.81	0.46
1:B:111:ASP:OD2	1:B:125:HIS:ND1	2.46	0.45
1:A:196:TRP:CH2	2:C:51:LYS:HB3	2.51	0.45
1:A:280:ALA:O	1:A:282:ASP:N	2.44	0.44
2:C:90:SER:O	2:C:93:THR:HG22	2.16	0.44
1:A:156:GLN:O	1:A:159:GLU:HG2	2.17	0.44
2:C:35:ILE:HG12	2:C:111:MET:HB2	1.98	0.44
2:D:24:ARG:NH1	2:D:30:LYS:O	2.50	0.44
1:B:283:LEU:O	1:B:286:MET:HG2	2.16	0.44
1:A:114:VAL:HG22	1:B:250:GLN:OE1	2.18	0.43
1:B:110:GLU:OE2	1:B:128:HIS:HA	2.18	0.43
2:D:65:LYS:HE2	2:D:65:LYS:HB3	1.85	0.43
1:A:124:GLU:HG3	1:A:125:HIS:CD2	2.54	0.42
2:D:11:ARG:HH21	2:D:15:GLN:HB3	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:ILE:HG21	1:B:209:GLU:HB2	2.02	0.42
1:A:219:SER:O	1:A:223:MET:HG2	2.19	0.42
2:D:51:LYS:HB3	1:B:196:TRP:CH2	2.55	0.42
1:A:133:MET:HE2	1:A:133:MET:HB3	1.86	0.41
1:A:206:GLU:OE1	2:C:70:ARG:NH2	2.54	0.41
1:A:196:TRP:CZ3	2:C:51:LYS:HB3	2.55	0.41
1:A:205:LEU:HD23	1:A:205:LEU:HA	1.87	0.41
1:A:226:GLN:O	1:A:230:MET:HG2	2.21	0.41
1:B:166:ARG:O	1:B:169:LYS:HG2	2.20	0.41
2:D:58:VAL:HG13	2:D:62:GLU:HB3	2.02	0.40
1:B:212:ILE:HG23	1:B:270:PHE:HE1	1.86	0.40
2:D:35:ILE:HB	2:D:52:PHE:HE2	1.86	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	191/193 (99%)	186 (97%)	5 (3%)	0	100	100
1	B	187/193 (97%)	183 (98%)	4 (2%)	0	100	100
2	C	112/114 (98%)	107 (96%)	5 (4%)	0	100	100
2	D	111/114 (97%)	106 (96%)	5 (4%)	0	100	100
All	All	601/614 (98%)	582 (97%)	19 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	193/193 (100%)	0.92	25 (12%) 7 7	59, 91, 133, 148	0
1	B	191/193 (98%)	0.86	24 (12%) 8 8	58, 86, 116, 141	0
2	C	114/114 (100%)	0.64	9 (7%) 18 19	37, 73, 115, 154	0
2	D	113/114 (99%)	0.64	9 (7%) 18 19	41, 82, 125, 163	0
All	All	611/614 (99%)	0.80	67 (10%) 10 10	37, 86, 125, 163	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	117	GLU	6.1
1	B	277	VAL	5.2
1	A	279	ARG	5.0
1	B	110	GLU	4.8
1	A	281	PRO	4.6
1	B	280	ALA	4.6
2	C	116	GLN	4.4
1	B	287	LEU	4.1
2	C	4	GLU	4.0
2	D	82	LEU	3.8
1	A	280	ALA	3.5
1	A	287	LEU	3.5
1	B	281	PRO	3.5
1	A	278	PHE	3.4
2	C	87	SER	3.3
1	B	117	TRP	3.2
1	A	277	VAL	3.2
1	A	276	ARG	3.1
1	A	284	LYS	3.1
2	D	89	VAL	3.1
2	D	87	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	151	LEU	3.0
2	C	89	VAL	3.0
1	A	135	GLU	2.9
1	A	182	ASN	2.9
2	C	80	PHE	2.9
1	B	124	GLU	2.9
1	A	230	MET	2.9
1	A	140	TYR	2.8
1	B	193	ILE	2.8
1	B	123	GLN	2.8
1	A	155	GLN	2.7
1	B	274	GLN	2.7
1	A	282	ASP	2.7
1	A	107	PHE	2.7
2	D	95	ILE	2.6
2	C	114	ALA	2.6
1	B	130	THR	2.5
2	D	38	TYR	2.5
1	A	123	GLN	2.5
1	B	230	MET	2.5
1	A	272	LYS	2.5
1	B	129	HIS	2.4
1	A	190	LEU	2.4
1	B	272	LYS	2.4
1	A	106	LEU	2.3
2	D	114	ALA	2.3
2	C	86	HIS	2.3
1	B	248	LEU	2.2
1	B	284	LYS	2.2
1	B	271	HIS	2.2
2	D	60	MET	2.2
1	B	115	ILE	2.2
1	B	99	ARG	2.1
1	B	190	LEU	2.1
1	B	282	ASP	2.1
1	B	109	GLN	2.1
1	B	203	GLN	2.1
2	D	85	GLY	2.1
1	A	121	ASP	2.1
1	A	275	ARG	2.1
1	A	172	TRP	2.1
1	A	283	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	273	ASN	2.1
1	B	136	VAL	2.1
2	D	113	TYR	2.0
2	C	78	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 oligosaccharides 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, 95th 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	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	A	301	1/1	0.97	0.08	108,108,108,108	0
3	ZN	A	302	1/1	0.97	0.05	114,114,114,114	0
3	ZN	B	301	1/1	0.98	0.05	75,75,75,75	0
3	ZN	B	302	1/1	0.98	0.05	84,84,84,84	0

6.5 Other polymers [i](#)

There are no such residues in this entry.