



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:26 AM UTC

PDB ID : 4TN3 / pdb_00004tn3
Title : Structure of the BBox-Coiled-coil region of Rhesus Trim5alpha
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Deposited on : 2014-06-03
Resolution : 3.20 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

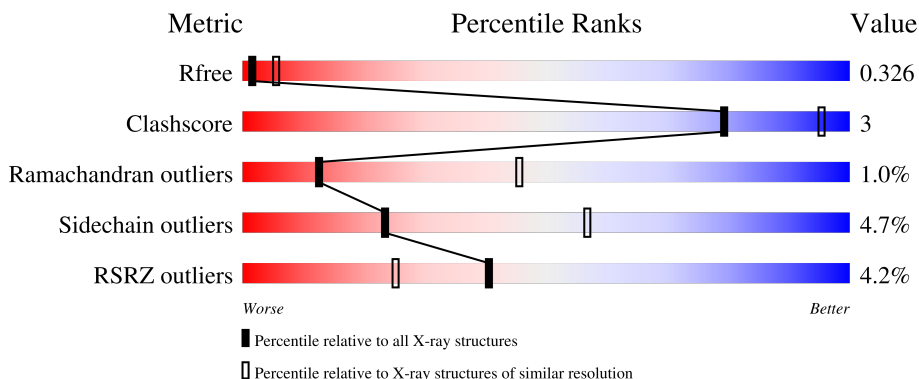
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	400	5% 79% 10% 10%
1	B	400	2% 80% 9% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11512 atoms, of which 5712 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 TRIM5/cyclophilin A fusion protein/T4 Lysozyme chimera.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	359	Total	C	H	N	O	S	0	0	0
			5740	1806	2854	508	554	18			
1	B	362	Total	C	H	N	O	S	0	0	0
			5768	1819	2858	518	557	16			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	68	MET	-	expression tag	UNP G9MAP5
A	69	ALA	-	expression tag	UNP G9MAP5
A	70	HIS	-	expression tag	UNP G9MAP5
A	71	HIS	-	expression tag	UNP G9MAP5
A	72	HIS	-	expression tag	UNP G9MAP5
A	73	HIS	-	expression tag	UNP G9MAP5
A	74	HIS	-	expression tag	UNP G9MAP5
A	75	HIS	-	expression tag	UNP G9MAP5
A	76	SER	-	expression tag	UNP G9MAP5
A	77	ALA	-	expression tag	UNP G9MAP5
A	78	ALA	-	expression tag	UNP G9MAP5
A	79	LEU	-	expression tag	UNP G9MAP5
A	80	GLU	-	expression tag	UNP G9MAP5
A	81	VAL	-	expression tag	UNP G9MAP5
A	82	LEU	-	expression tag	UNP G9MAP5
A	83	PHE	-	expression tag	UNP G9MAP5
A	84	GLN	-	expression tag	UNP G9MAP5
A	85	GLY	-	expression tag	UNP G9MAP5
A	86	PRO	-	expression tag	UNP G9MAP5
A	87	GLY	-	expression tag	UNP G9MAP5
A	315	GLY	ARG	conflict	UNP P00720
A	323	ASN	ASP	conflict	UNP P00720
A	357	THR	CYS	conflict	UNP P00720
A	400	ALA	CYS	conflict	UNP P00720
A	440	ARG	ILE	conflict	UNP P00720

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Chain	Residue	Modelled	Actual	Comment	Reference
B	68	MET	-	expression tag	UNP G9MAP5
B	69	ALA	-	expression tag	UNP G9MAP5
B	70	HIS	-	expression tag	UNP G9MAP5
B	71	HIS	-	expression tag	UNP G9MAP5
B	72	HIS	-	expression tag	UNP G9MAP5
B	73	HIS	-	expression tag	UNP G9MAP5
B	74	HIS	-	expression tag	UNP G9MAP5
B	75	HIS	-	expression tag	UNP G9MAP5
B	76	SER	-	expression tag	UNP G9MAP5
B	77	ALA	-	expression tag	UNP G9MAP5
B	78	ALA	-	expression tag	UNP G9MAP5
B	79	LEU	-	expression tag	UNP G9MAP5
B	80	GLU	-	expression tag	UNP G9MAP5
B	81	VAL	-	expression tag	UNP G9MAP5
B	82	LEU	-	expression tag	UNP G9MAP5
B	83	PHE	-	expression tag	UNP G9MAP5
B	84	GLN	-	expression tag	UNP G9MAP5
B	85	GLY	-	expression tag	UNP G9MAP5
B	86	PRO	-	expression tag	UNP G9MAP5
B	87	GLY	-	expression tag	UNP G9MAP5
B	315	GLY	ARG	conflict	UNP P00720
B	323	ASN	ASP	conflict	UNP P00720
B	357	THR	CYS	conflict	UNP P00720
B	400	ALA	CYS	conflict	UNP P00720
B	440	ARG	ILE	conflict	UNP P00720

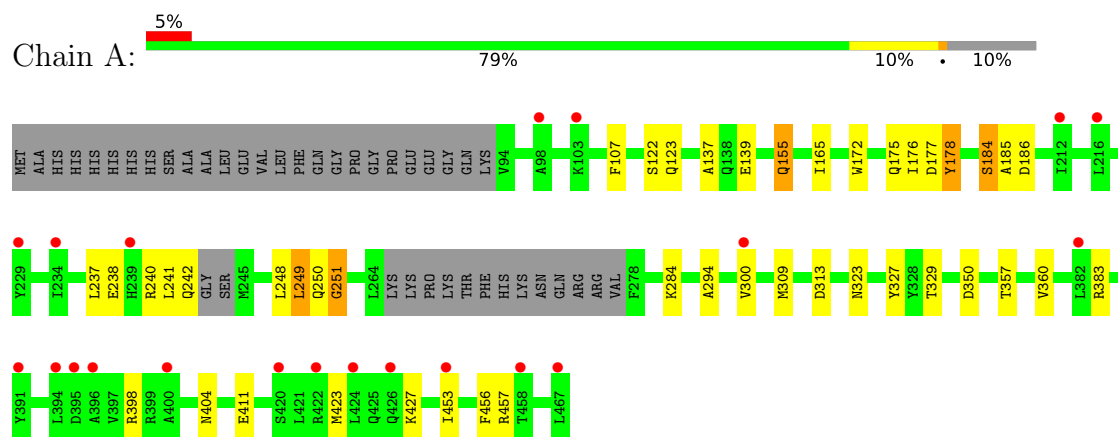
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

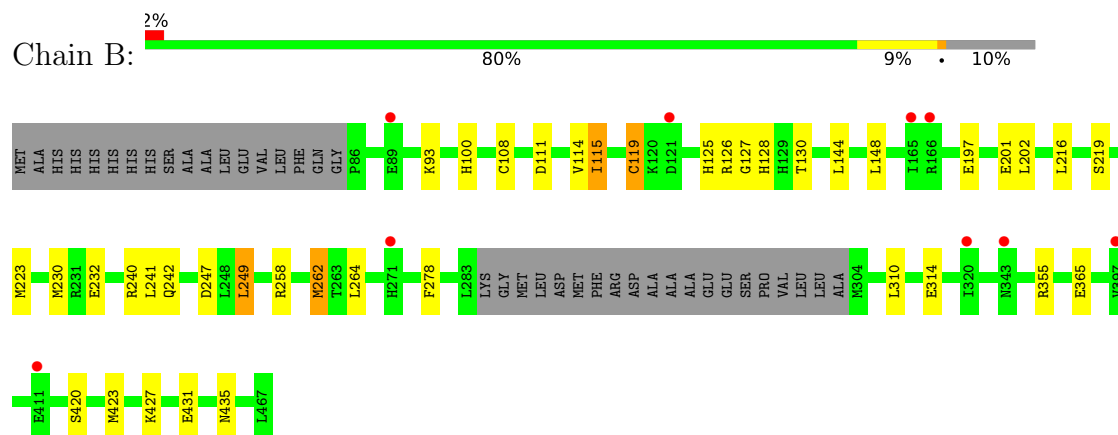
3 Residue-property plots

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: TRIM5/cyclophilin A fusion protein/T4 Lysozyme chimera



- Molecule 1: TRIM5/cyclophilin A fusion protein/T4 Lysozyme chimera



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	124.12Å 59.74Å 146.96Å 90.00° 94.77° 90.00°	Depositor
Resolution (Å)	33.93 – 3.20 33.93 – 3.20	Depositor EDS
% Data completeness (in resolution range)	98.9 (33.93-3.20) 98.8 (33.93-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 3.18Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.258 , 0.316 0.272 , 0.326	Depositor DCC
R_{free} test set	903 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	79.7	Xtriage
Anisotropy	0.463	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 80.3	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	0.91	EDS
Total number of atoms	11512	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.16% 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

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	0.33	0/2926	0.66	0/3933
1	B	0.33	0/2955	0.66	0/3974
All	All	0.33	0/5881	0.66	0/7907

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	2886	2854	2861	17	0
1	B	2910	2858	2867	16	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
All	All	5800	5712	5728	31	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 (31) 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:144:LEU:O	1:B:148:LEU:N	2.30	0.64
1:A:155:GLN:O	1:A:155:GLN:NE2	2.33	0.62
1:B:125:HIS:O	1:B:127:GLY:N	2.33	0.61
1:A:249:LEU:O	1:A:251:GLY:N	2.37	0.58
1:A:423:MET:O	1:A:427:LYS:N	2.38	0.56
1:B:431:GLU:N	1:B:431:GLU:OE1	2.39	0.56
1:A:284:LYS:NZ	1:B:197:GLU:OE2	2.36	0.55
1:A:323:ASN:N	1:A:327:TYR:O	2.39	0.55
1:B:423:MET:O	1:B:427:LYS:N	2.42	0.52
1:B:115:ILE:HB	1:B:119:CYS:HB3	1.93	0.51
1:B:240:ARG:NH1	1:B:247:ASP:O	2.44	0.50
1:B:310:LEU:O	1:B:314:GLU:N	2.42	0.50
1:A:176:ILE:C	1:A:178:TYR:H	2.19	0.49
1:B:230:MET:HE2	1:B:262:MET:SD	2.53	0.48
1:A:350:ASP:OD1	1:A:357:THR:N	2.43	0.47
1:A:238:GLU:O	1:A:242:GLN:NE2	2.49	0.46
1:A:453:ILE:O	1:A:456:PHE:N	2.49	0.45
1:A:184:SER:OG	1:A:185:ALA:N	2.49	0.45
1:A:309:MET:HE2	1:A:404:ASN:HD22	1.82	0.45
1:B:355:ARG:NH2	1:B:365:GLU:OE1	2.48	0.44
1:A:398:ARG:NH2	1:A:457:ARG:O	2.47	0.44
1:B:108:CYS:SG	1:B:111:ASP:HB2	2.57	0.44
1:A:107:PHE:CD1	1:B:249:LEU:HD22	2.53	0.44
1:A:237:LEU:HA	1:A:240:ARG:HB3	1.98	0.44
1:A:172:TRP:HA	1:A:175:GLN:HB3	2.00	0.43
1:A:383:ARG:NH2	1:A:411:GLU:OE2	2.52	0.43
1:B:111:ASP:OD2	1:B:128:HIS:CE1	2.71	0.43
1:B:420:SER:OG	1:B:435:ASN:OD1	2.34	0.43
1:B:100:HIS:CE1	1:B:119:CYS:HA	2.54	0.42
1:A:122:SER:OG	1:A:123:GLN:N	2.54	0.41
1:B:240:ARG:C	1:B:242:GLN:H	2.29	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	353/400 (88%)	312 (88%)	35 (10%)	6 (2%)	7	35
1	B	358/400 (90%)	331 (92%)	26 (7%)	1 (0%)	36	68
All	All	711/800 (89%)	643 (90%)	61 (9%)	7 (1%)	12	45

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	250	GLN
1	B	126	ARG
1	A	137	ALA
1	A	241	LEU
1	A	294	ALA
1	A	177	ASP
1	A	251	GLY

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	308/349 (88%)	296 (96%)	12 (4%)	28	62
1	B	310/349 (89%)	293 (94%)	17 (6%)	19	52
All	All	618/698 (88%)	589 (95%)	29 (5%)	23	57

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	139	GLU
1	A	155	GLN
1	A	165	ILE
1	A	178	TYR
1	A	184	SER
1	A	186	ASP

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Mol	Chain	Res	Type
1	A	248	LEU
1	A	249	LEU
1	A	300	VAL
1	A	313	ASP
1	A	329	THR
1	A	360	VAL
1	B	93	LYS
1	B	114	VAL
1	B	115	ILE
1	B	119	CYS
1	B	130	THR
1	B	201	GLU
1	B	202	LEU
1	B	216	LEU
1	B	219	SER
1	B	223	MET
1	B	232	GLU
1	B	241	LEU
1	B	249	LEU
1	B	258	ARG
1	B	262	MET
1	B	264	LEU
1	B	278	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	GLN
1	A	156	GLN
1	A	384	ASN
1	A	404	ASN
1	B	145	GLN
1	B	426	GLN

5.3.3 RNA ⓘ

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	359/400 (89%)	0.24	21 (5%) 29 18	16, 86, 170, 233	0
1	B	362/400 (90%)	0.35	9 (2%) 58 39	32, 108, 185, 270	0
All	All	721/800 (90%)	0.29	30 (4%) 40 25	16, 97, 182, 270	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	165	ILE	5.0
1	A	426	GLN	3.2
1	A	216	LEU	3.0
1	A	234	ILE	3.0
1	B	397	VAL	2.9
1	B	320	ILE	2.9
1	B	89	GLU	2.9
1	A	422	ARG	2.9
1	A	400	ALA	2.8
1	A	420	SER	2.6
1	A	229	TYR	2.6
1	A	239	HIS	2.6
1	B	166	ARG	2.6
1	A	424	LEU	2.5
1	A	382	LEU	2.5
1	A	98	ALA	2.3
1	A	458	THR	2.3
1	A	394	LEU	2.3
1	A	396	ALA	2.2
1	A	453	ILE	2.1
1	A	103	LYS	2.1
1	A	212	ILE	2.1
1	A	391	TYR	2.1
1	B	411	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	300	VAL	2.1
1	A	395	ASP	2.1
1	A	467	LEU	2.0
1	B	343	ASN	2.0
1	B	271	HIS	2.0
1	B	121	ASP	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(Å ²)	Q<0.9
2	ZN	A	701	1/1	0.98	0.03	76,76,76,76	0
2	ZN	B	700	1/1	0.98	0.03	66,66,66,66	0
2	ZN	A	702	1/1	0.99	0.02	68,68,68,68	0
2	ZN	B	701	1/1	0.99	0.03	57,57,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.