



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:48 AM UTC

PDB ID : 3QK9 / pdb_00003qk9
Title : Yeast Tim44 C-terminal domain complexed with Cymal-3
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Deposited on : 2011-01-31
Resolution : 3.10 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

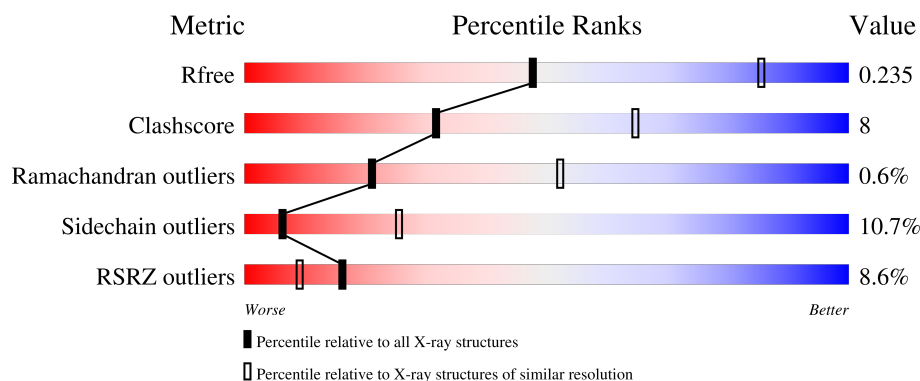
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	222	8% 55% 20% • 22%
1	B	222	6% 57% 19% • 21%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2845 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 Mitochondrial import inner membrane translocase subunit TIM44.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	174	Total	C	N	O	S	0	0	0
			1414	908	242	260	4			
1	B	175	Total	C	N	O	S	0	0	0
			1422	914	243	261	4			

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cl	0	0
			1	1		
2	B	1	Total	Cl	0	0
			1	1		

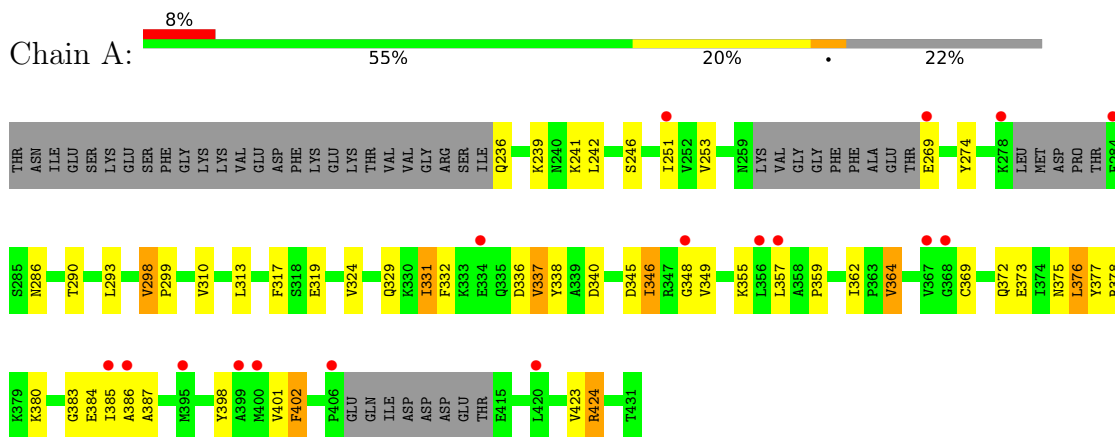
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total	O	0	0
			7	7		

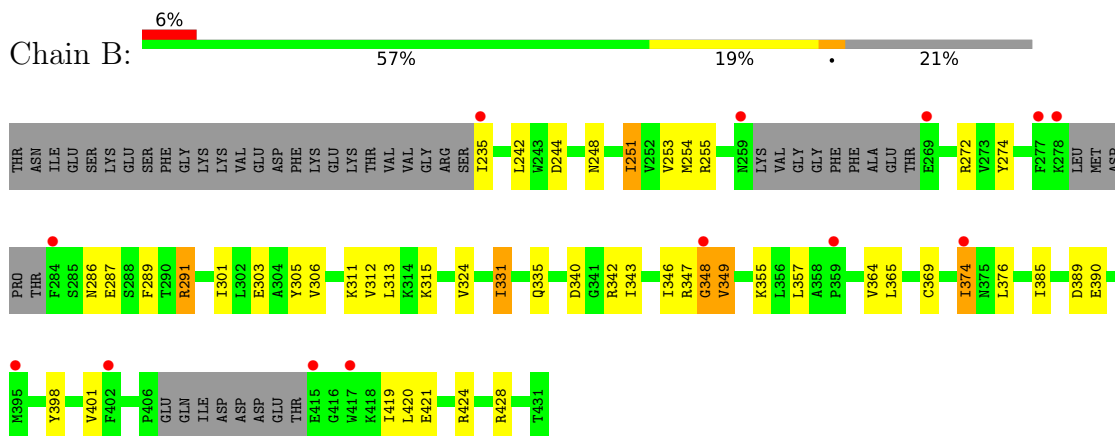
3 Residue-property plots [i](#)

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: Mitochondrial import inner membrane translocase subunit TIM44



- Molecule 1: Mitochondrial import inner membrane translocase subunit TIM44



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	100.59Å 100.59Å 209.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.93 – 3.10 32.93 – 3.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (32.93-3.10) 93.8 (32.93-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.235 , 0.317 0.313 , 0.235	Depositor DCC
R_{free} test set	691 reflections (6.10%)	wwPDB-VP
Wilson B-factor (Å ²)	95.2	Xtriage
Anisotropy	0.221	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 51.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	2845	wwPDB-VP
Average B, all atoms (Å ²)	106.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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: CL

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	0.53	1/1439 (0.1%)	0.87	5/1933 (0.3%)
1	B	0.45	0/1447	0.88	4/1944 (0.2%)
All	All	0.49	1/2886 (0.0%)	0.87	9/3877 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	298	VAL	CA-CB	6.27	1.57	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	349	VAL	N-CA-CB	8.87	125.86	111.23
1	B	348	GLY	N-CA-C	-8.55	92.91	113.18
1	A	387	ALA	N-CA-C	8.07	122.60	109.85
1	B	390	GLU	N-CA-C	-6.93	104.45	113.12
1	A	387	ALA	N-CA-CB	-6.23	101.11	111.20
1	A	402	PHE	N-CA-C	5.38	118.84	110.17
1	B	389	ASP	CB-CA-C	5.21	119.78	109.66
1	A	386	ALA	CB-CA-C	5.20	120.76	110.42
1	A	386	ALA	N-CA-C	-5.01	100.12	110.80

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	1414	0	1429	25	0
1	B	1422	0	1440	24	0
2	A	1	0	0	0	0
2	B	1	0	0	1	0
3	A	7	0	0	0	0
All	All	2845	0	2869	46	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 8.

All (46) 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:348:GLY:O	1:B:369:CYS:HB2	1.74	0.85
1:A:246:SER:HB3	1:A:251:ILE:HD11	1.69	0.74
1:B:244:ASP:HA	1:B:255:ARG:HE	1.57	0.68
1:A:329:GLN:HA	1:A:332:PHE:HD1	1.58	0.68
1:A:331:ILE:HD11	1:B:331:ILE:HD13	1.75	0.67
1:A:375:ASN:HD21	1:B:335:GLN:HE22	1.44	0.65
1:A:348:GLY:O	1:A:369:CYS:HA	1.98	0.64
1:A:313:LEU:O	1:A:317:PHE:HB2	2.00	0.61
1:A:329:GLN:HA	1:A:332:PHE:CD1	2.37	0.60
1:B:348:GLY:O	1:B:369:CYS:CB	2.50	0.58
1:A:364:VAL:HG13	1:A:401:VAL:HG13	1.86	0.57
1:A:369:CYS:HB3	1:A:398:TYR:HB2	1.88	0.56
1:A:345:ASP:HB3	1:A:372:GLN:HB2	1.90	0.54
1:B:301:ILE:HD13	1:B:419:ILE:HD11	1.88	0.54
1:B:342:ARG:HG2	1:B:374:ILE:HD11	1.89	0.54
1:B:342:ARG:HG2	1:B:374:ILE:CD1	2.39	0.53
1:B:251:ILE:HA	1:B:254:MET:HB2	1.92	0.52
1:B:343:ILE:HD12	1:B:346:ILE:HD11	1.92	0.51
1:A:324:VAL:HA	1:B:324:VAL:HG23	1.92	0.51
1:A:253:VAL:HG23	1:A:423:VAL:HG22	1.94	0.50
1:B:274:TYR:HE2	1:B:286:ASN:HB2	1.76	0.50
1:B:398:TYR:CE2	1:B:424:ARG:HG3	2.47	0.50
1:B:303:GLU:HB3	1:B:343:ILE:HD11	1.94	0.50
1:A:357:LEU:HB3	1:A:359:PRO:HD2	1.93	0.50
1:B:244:ASP:HA	1:B:255:ARG:NE	2.26	0.49
1:B:398:TYR:CZ	1:B:424:ARG:HG3	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:428:ARG:NH1	2:B:1:CL:CL	2.81	0.49
1:A:290:THR:HA	1:A:293:LEU:HD12	1.98	0.46
1:A:298:VAL:HB	1:A:299:PRO:HD3	1.99	0.45
1:B:305:TYR:HA	1:B:313:LEU:HD11	1.97	0.45
1:B:248:ASN:O	1:B:251:ILE:HD13	2.16	0.45
1:A:337:VAL:HG22	1:A:377:TYR:HB3	1.99	0.44
1:B:401:VAL:HB	1:B:421:GLU:H	1.84	0.43
1:B:306:VAL:HG21	1:B:343:ILE:HG23	2.01	0.42
1:B:312:VAL:HA	1:B:315:LYS:HE2	2.00	0.42
1:A:346:ILE:HD11	1:A:349:VAL:HG21	2.00	0.42
1:A:401:VAL:O	1:A:401:VAL:HG12	2.19	0.42
1:B:287:GLU:O	1:B:291:ARG:HD2	2.19	0.42
1:A:338:TYR:CE1	1:A:378:ARG:HB2	2.55	0.42
1:A:274:TYR:HE2	1:A:286:ASN:HD22	1.68	0.41
1:A:373:GLU:OE1	1:A:424:ARG:NH1	2.54	0.41
1:A:340:ASP:O	1:A:376:LEU:HB3	2.20	0.41
1:A:378:ARG:NE	1:A:384:GLU:O	2.53	0.41
1:B:340:ASP:O	1:B:376:LEU:O	2.38	0.41
1:A:336:ASP:HB3	1:A:380:LYS:HG2	2.03	0.41
1:A:364:VAL:HA	1:A:402:PHE:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	166/222 (75%)	152 (92%)	13 (8%)	1 (1%)	21	52
1	B	167/222 (75%)	158 (95%)	8 (5%)	1 (1%)	21	52
All	All	333/444 (75%)	310 (93%)	21 (6%)	2 (1%)	21	52

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	349	VAL
1	A	383	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	153/196 (78%)	137 (90%)	16 (10%)	6	26
1	B	154/196 (79%)	137 (89%)	17 (11%)	6	24
All	All	307/392 (78%)	274 (89%)	33 (11%)	6	25

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	236	GLN
1	A	239	LYS
1	A	241	LYS
1	A	242	LEU
1	A	269	GLU
1	A	310	VAL
1	A	319	GLU
1	A	331	ILE
1	A	337	VAL
1	A	346	ILE
1	A	355	LYS
1	A	362	ILE
1	A	364	VAL
1	A	376	LEU
1	A	385	ILE
1	A	424	ARG
1	B	235	ILE
1	B	242	LEU
1	B	251	ILE
1	B	253	VAL
1	B	272	ARG
1	B	289	PHE

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Mol	Chain	Res	Type
1	B	291	ARG
1	B	311	LYS
1	B	331	ILE
1	B	347	ARG
1	B	355	LYS
1	B	357	LEU
1	B	364	VAL
1	B	365	LEU
1	B	374	ILE
1	B	385	ILE
1	B	420	LEU

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	323	ASN
1	A	375	ASN
1	B	236	GLN
1	B	276	GLN
1	B	323	ASN
1	B	329	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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	174/222 (78%)	0.78	17 (9%) 13 7	78, 104, 145, 162	0
1	B	175/222 (78%)	0.62	13 (7%) 20 11	74, 100, 148, 162	0
All	All	349/444 (78%)	0.70	30 (8%) 16 9	74, 102, 147, 162	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	368	GLY	5.1
1	A	357	LEU	4.3
1	B	395	MET	4.2
1	A	284	PHE	3.9
1	A	395	MET	3.6
1	B	278	LYS	3.5
1	A	400	MET	3.3
1	B	235	ILE	3.3
1	B	348	GLY	3.3
1	B	402	PHE	3.2
1	B	417	TRP	3.0
1	B	415	GLU	3.0
1	A	334	GLU	2.8
1	B	284	PHE	2.7
1	B	359	PRO	2.6
1	A	367	VAL	2.6
1	A	399	ALA	2.6
1	A	278	LYS	2.5
1	A	386	ALA	2.5
1	B	374	ILE	2.5
1	A	356	LEU	2.5
1	A	269	GLU	2.5
1	B	269	GLU	2.5
1	A	420	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	259	ASN	2.2
1	A	406	PRO	2.2
1	A	385	ILE	2.2
1	B	277	PHE	2.2
1	A	348	GLY	2.1
1	A	251	ILE	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
2	CL	B	1	1/1	0.81	0.13	48,48,48,48	0
2	CL	A	1	1/1	0.88	0.12	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.