



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:55 AM UTC

PDB ID : 3IVF / pdb_00003ivf
Title : Crystal structure of the talin head FERM domain
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Deposited on : 2009-09-01
Resolution : 1.94 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

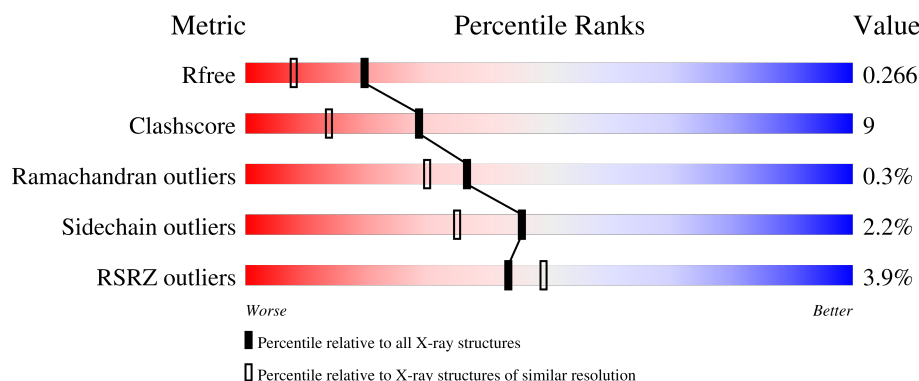
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	371	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3455 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 Talin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	358	Total	C	N	O	S	42	0	0
			2899	1844	501	536	18			

There are 29 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLU	deletion	UNP P26039
A	?	-	GLY	deletion	UNP P26039
A	?	-	THR	deletion	UNP P26039
A	?	-	GLY	deletion	UNP P26039
A	?	-	THR	deletion	UNP P26039
A	?	-	LEU	deletion	UNP P26039
A	?	-	ARG	deletion	UNP P26039
A	?	-	LYS	deletion	UNP P26039
A	?	-	ASP	deletion	UNP P26039
A	?	-	LYS	deletion	UNP P26039
A	?	-	THR	deletion	UNP P26039
A	?	-	LEU	deletion	UNP P26039
A	?	-	LEU	deletion	UNP P26039
A	?	-	ARG	deletion	UNP P26039
A	?	-	ASP	deletion	UNP P26039
A	?	-	GLU	deletion	UNP P26039
A	?	-	LYS	deletion	UNP P26039
A	?	-	LYS	deletion	UNP P26039
A	?	-	MET	deletion	UNP P26039
A	?	-	GLU	deletion	UNP P26039
A	?	-	LYS	deletion	UNP P26039
A	?	-	LEU	deletion	UNP P26039
A	?	-	LYS	deletion	UNP P26039
A	?	-	GLN	deletion	UNP P26039
A	?	-	LYS	deletion	UNP P26039
A	?	-	LEU	deletion	UNP P26039
A	?	-	HIS	deletion	UNP P26039

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	THR	deletion	UNP P26039
A	?	-	ASP	deletion	UNP P26039

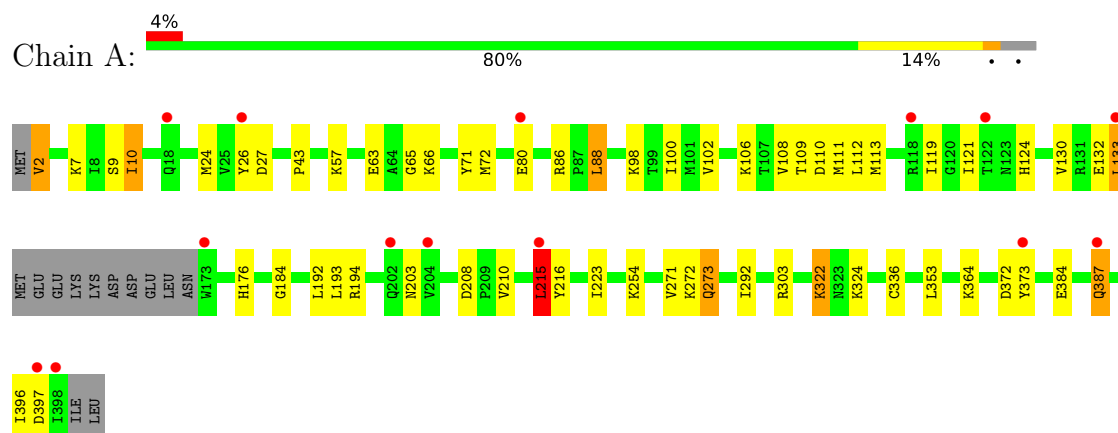
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	556	Total	O	0	0
			556	556		

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: Talin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	40.26Å 72.16Å 162.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.38 – 1.94 43.38 – 1.94	Depositor EDS
% Data completeness (in resolution range)	96.1 (43.38-1.94) 96.1 (43.38-1.94)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.193 , 0.257 0.206 , 0.266	Depositor DCC
R_{free} test set	1721 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3455	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.00% 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	1.37	8/2954 (0.3%)	1.13	7/3971 (0.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	2	VAL	C-N	-9.87	1.20	1.33
1	A	88	LEU	CA-C	5.92	1.60	1.52
1	A	322	LYS	CG-CD	-5.77	1.35	1.52
1	A	215	LEU	C-O	-5.54	1.17	1.24
1	A	216	TYR	C-O	-5.41	1.17	1.24
1	A	194	ARG	C-O	-5.28	1.18	1.24
1	A	223	ILE	CA-CB	-5.08	1.48	1.54
1	A	130	VAL	N-CA	-5.02	1.40	1.46

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	ILE	CB-CA-C	-6.03	103.54	110.41
1	A	43	PRO	O-C-N	5.64	123.90	121.31
1	A	210	VAL	CB-CA-C	-5.59	104.69	112.02
1	A	324	LYS	CA-CB-CG	-5.15	103.80	114.10
1	A	336	CYS	N-CA-C	5.14	116.10	108.60
1	A	208	ASP	CA-C-N	5.09	124.70	119.05
1	A	208	ASP	C-N-CA	5.09	124.70	119.05

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	2899	0	2921	50	1
2	A	556	0	0	16	1
All	All	3455	0	2921	50	1

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 9.

All (50) 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:303:ARG:NH1	2:A:423:HOH:O	2.00	0.92
1:A:112:LEU:HD23	2:A:871:HOH:O	1.80	0.82
1:A:384:GLU:O	1:A:387:GLN:HG2	1.81	0.80
1:A:373:TYR:CD2	2:A:410:HOH:O	2.35	0.80
1:A:110:ASP:OD2	2:A:541:HOH:O	2.02	0.78
1:A:88:LEU:HD22	1:A:111:MET:HE1	1.65	0.77
1:A:203:ASN:ND2	2:A:432:HOH:O	2.12	0.77
1:A:373:TYR:HD2	2:A:410:HOH:O	1.69	0.73
1:A:2:VAL:HG22	1:A:2:VAL:O	1.89	0.72
1:A:215:LEU:O	1:A:215:LEU:HD12	1.93	0.69
1:A:7:LYS:NZ	2:A:755:HOH:O	2.26	0.67
1:A:88:LEU:CD2	1:A:111:MET:HE1	2.25	0.67
1:A:387:GLN:NE2	2:A:430:HOH:O	2.26	0.67
1:A:24:MET:HE1	1:A:65:GLY:O	1.95	0.66
1:A:271:VAL:O	1:A:273:GLN:N	2.29	0.64
1:A:119:ILE:HD11	1:A:121:ILE:HD12	1.81	0.61
1:A:102:VAL:HG11	1:A:111:MET:HE3	1.84	0.59
1:A:2:VAL:O	1:A:2:VAL:CG2	2.51	0.57
1:A:271:VAL:C	1:A:273:GLN:H	2.13	0.56
1:A:102:VAL:CB	1:A:111:MET:HE3	2.35	0.56
1:A:63:GLU:HB2	1:A:66:LYS:HG2	1.89	0.55
1:A:373:TYR:CE2	2:A:410:HOH:O	2.59	0.55
1:A:57:LYS:HE2	1:A:184:GLY:HA2	1.91	0.53
1:A:102:VAL:HG21	1:A:111:MET:HE3	1.92	0.51
1:A:57:LYS:HE2	1:A:184:GLY:CA	2.40	0.51
1:A:9:SER:HB3	1:A:80:GLU:HG2	1.94	0.50
1:A:303:ARG:NH2	2:A:453:HOH:O	2.36	0.50
1:A:102:VAL:CG1	1:A:111:MET:HE3	2.42	0.49
1:A:108:VAL:HB	1:A:176:HIS:HA	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:TYR:C	1:A:72:MET:HG2	2.40	0.46
1:A:396:ILE:O	1:A:397:ASP:C	2.59	0.46
1:A:109:THR:OG1	1:A:176:HIS:HB3	2.16	0.45
1:A:192:LEU:HG	1:A:193:LEU:N	2.31	0.45
1:A:303:ARG:NE	2:A:453:HOH:O	2.49	0.44
1:A:26:TYR:CE1	1:A:27:ASP:OD1	2.71	0.43
1:A:133:LEU:HD12	1:A:133:LEU:HA	1.40	0.43
1:A:106:LYS:HD2	2:A:541:HOH:O	2.17	0.43
1:A:86:ARG:NH1	2:A:865:HOH:O	2.43	0.43
1:A:102:VAL:HG21	1:A:111:MET:CE	2.48	0.42
1:A:254:LYS:HE3	1:A:254:LYS:HB2	1.80	0.42
1:A:292:ILE:HD13	2:A:590:HOH:O	2.19	0.42
1:A:98:LYS:HB3	1:A:100:ILE:HD11	2.02	0.42
1:A:26:TYR:C	1:A:26:TYR:CD1	2.97	0.42
1:A:102:VAL:CG2	1:A:111:MET:HE3	2.50	0.41
1:A:364:LYS:HE3	2:A:497:HOH:O	2.20	0.41
1:A:113:MET:HG2	1:A:124:HIS:CE1	2.55	0.41
1:A:111:MET:HE2	1:A:111:MET:HB3	1.79	0.40
1:A:102:VAL:HB	1:A:111:MET:HE3	2.03	0.40
1:A:271:VAL:C	1:A:273:GLN:N	2.73	0.40
1:A:372:ASP:OD1	2:A:866:HOH:O	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:LYS:NZ	2:A:925:HOH:O[4_454]	2.05	0.15

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	354/371 (95%)	350 (99%)	3 (1%)	1 (0%)	36	30

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	272	LYS

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	317/330 (96%)	310 (98%)	7 (2%)	45	35

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ILE
1	A	132	GLU
1	A	133	LEU
1	A	215	LEU
1	A	273	GLN
1	A	353	LEU
1	A	387	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	242	GLN
1	A	246	GLN
1	A	281	GLN
1	A	288	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](#)

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	358/371 (96%)	0.26	14 (3%)	43 49	12, 27, 44, 62	10 (2%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	373	TYR	3.8
1	A	173	TRP	3.2
1	A	215	LEU	2.9
1	A	26	TYR	2.7
1	A	202	GLN	2.7
1	A	133	LEU	2.5
1	A	118	ARG	2.5
1	A	397	ASP	2.4
1	A	204	VAL	2.4
1	A	80	GLU	2.3
1	A	398	ILE	2.3
1	A	122	THR	2.2
1	A	387	GLN	2.1
1	A	18	GLN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.