



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 9QN7 / pdb_00009qn7
Title : Structure of talin in complex with a peptide fragment
Authors : Zacharchenko, T.
Deposited on : 2025-03-24
Resolution : 2.76 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
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)	:	2.46

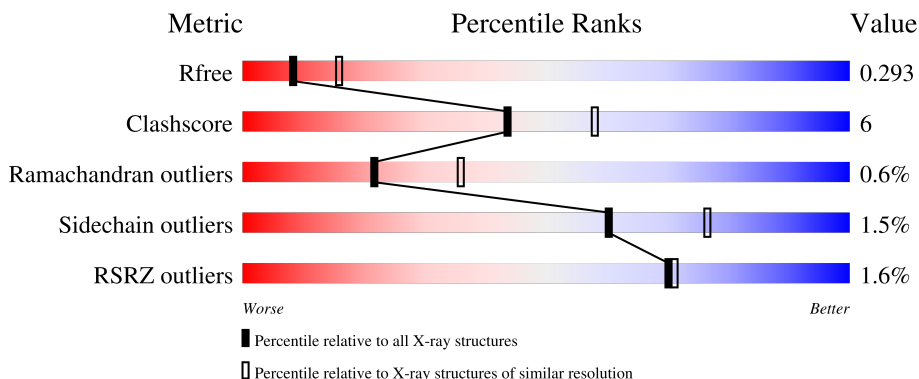
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 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}	164625	1606 (2.78-2.74)
Clashscore	180529	1689 (2.78-2.74)
Ramachandran outliers	177936	1665 (2.78-2.74)
Sidechain outliers	177891	1665 (2.78-2.74)
RSRZ outliers	164620	1606 (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 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	330	<div> <div>0%</div> <div>83%</div> <div>15%</div> <div>•</div> </div>
1	B	330	<div> <div>2%</div> <div>83%</div> <div>14%</div> <div>• •</div> </div>
2	C	27	<div> <div>59%</div> <div>30%</div> <div>11%</div> </div>
2	D	27	<div> <div>7%</div> <div>33%</div> <div>15%</div> <div>•</div> <div>48%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5102 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	325	Total	C	N	O	S	0	0	0
			2395	1487	424	473	11			
1	B	321	Total	C	N	O	S	0	1	0
			2373	1475	420	466	12			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1964	GLY	-	expression tag	UNP P26039
A	1965	THR	-	expression tag	UNP P26039
A	1966	PHE	-	expression tag	UNP P26039
A	1967	THR	-	expression tag	UNP P26039
A	1968	GLY	-	expression tag	UNP P26039
A	1969	ILE	-	expression tag	UNP P26039
A	1970	ASP	-	expression tag	UNP P26039
A	1971	PRO	-	expression tag	UNP P26039
A	1972	PHE	-	expression tag	UNP P26039
A	1973	THR	-	expression tag	UNP P26039
B	1964	GLY	-	expression tag	UNP P26039
B	1965	THR	-	expression tag	UNP P26039
B	1966	PHE	-	expression tag	UNP P26039
B	1967	THR	-	expression tag	UNP P26039
B	1968	GLY	-	expression tag	UNP P26039
B	1969	ILE	-	expression tag	UNP P26039
B	1970	ASP	-	expression tag	UNP P26039
B	1971	PRO	-	expression tag	UNP P26039
B	1972	PHE	-	expression tag	UNP P26039
B	1973	THR	-	expression tag	UNP P26039

- Molecule 2 is a protein called Tensin-3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	24	Total	C	N	O	0	0	0
			197	125	30	42			
2	D	14	Total	C	N	O	0	0	0
			116	72	20	24			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	10	Total	O	0	0
			10	10		
3	B	9	Total	O	0	0
			9	9		
3	C	2	Total	O	0	0
			2	2		

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	84.94Å 84.94Å 346.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	76.26 – 2.76 76.26 – 2.76	Depositor EDS
% Data completeness (in resolution range)	99.6 (76.26-2.76) 99.6 (76.26-2.76)	Depositor EDS
R_{merge}	0.44	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.257 , 0.296 0.258 , 0.293	Depositor DCC
R_{free} test set	1692 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	64.3	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 86.2	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	0.95	EDS
Total number of atoms	5102	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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 [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	0.24	0/2421	0.61	0/3285
1	B	0.22	0/2402	0.60	0/3257
2	C	0.21	0/199	0.31	0/271
2	D	0.13	0/115	0.34	0/154
All	All	0.23	0/5137	0.59	0/6967

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	2395	0	2460	31	0
1	B	2373	0	2445	28	0
2	C	197	0	194	8	0
2	D	116	0	119	4	0
3	A	10	0	0	0	0
3	B	9	0	0	1	0
3	C	2	0	0	0	0
All	All	5102	0	5218	62	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 6.

The worst 5 of 62 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:2222:ALA:HB1	1:B:2290:MET:HE1	1.66	0.78
1:B:2198:GLN:NE2	3:B:2301:HOH:O	2.24	0.70
1:B:2037:ALA:HA	1:B:2044:LEU:HD12	1.80	0.64
1:A:1993:ASP:OD1	1:A:2085:LYS:NZ	2.27	0.60
2:D:694:ASP:OD1	2:D:694:ASP:N	2.34	0.60

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	323/330 (98%)	314 (97%)	6 (2%)	3 (1%)	14	26
1	B	320/330 (97%)	315 (98%)	4 (1%)	1 (0%)	37	55
2	C	22/27 (82%)	21 (96%)	1 (4%)	0	100	100
2	D	12/27 (44%)	10 (83%)	2 (17%)	0	100	100
All	All	677/714 (95%)	660 (98%)	13 (2%)	4 (1%)	22	36

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2106	GLY
1	A	2008	GLY
1	B	2106	GLY
1	A	1971	PRO

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	252/255 (99%)	250 (99%)	2 (1%)	79	88
1	B	250/255 (98%)	244 (98%)	6 (2%)	44	65
2	C	24/27 (89%)	24 (100%)	0	100	100
2	D	14/27 (52%)	13 (93%)	1 (7%)	12	22
All	All	540/564 (96%)	531 (98%)	9 (2%)	60	74

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	2288	GLU
2	D	694	ASP
1	B	2170	SER
1	B	2184[A]	MET
1	B	2184[B]	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2049	GLN
1	B	2041	GLN
1	B	2082	ASN
1	B	2152	HIS
1	B	2262	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

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	325/330 (98%)	-0.22	4 (1%) 76 78	51, 78, 126, 195	0
1	B	321/330 (97%)	-0.13	5 (1%) 70 71	51, 83, 133, 165	1 (0%)
2	C	24/27 (88%)	-0.19	0 100 100	75, 100, 136, 139	0
2	D	14/27 (51%)	0.72	2 (14%) 7 9	128, 148, 182, 192	0
All	All	684/714 (95%)	-0.16	11 (1%) 70 71	51, 82, 135, 195	1 (0%)

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1972	PHE	3.8
1	B	1971	PRO	3.8
1	B	2105	VAL	3.4
2	D	695	ILE	3.1
2	D	707	LEU	3.1

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.