



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 21, 2021 – 08:02 PM BST

PDB ID : 7A8T  
Title : Crystal structure of sarcomeric protein FATZ-1 (mini-FATZ-1 construct) in complex with rod domain of alpha-actinin-2  
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Deposited on : 2020-08-31  
Resolution : 2.69 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](mailto: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.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.20  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.20

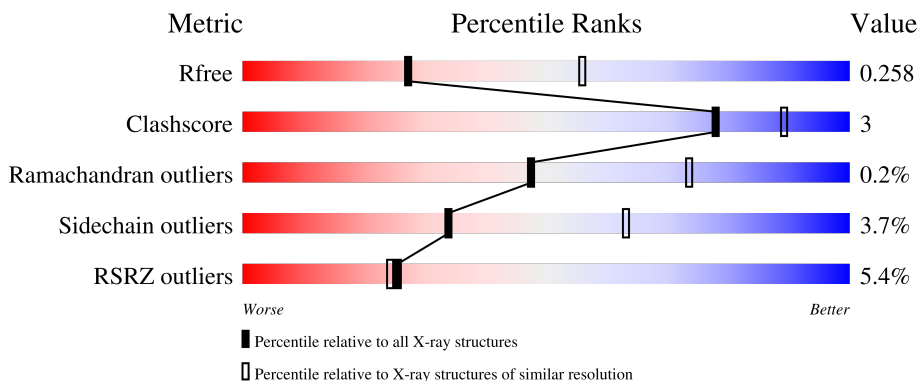
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.69 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))
$R_{free}$	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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  $\geq 3$ , 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	476	 5% 87% 12%
2	B	92	 7% 47% 51%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4309 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 Alpha-actinin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	474	3935	2450	713	755	17	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	271	GLY	-	expression tag	UNP P35609
A	272	SER	-	expression tag	UNP P35609
A	273	SER	-	expression tag	UNP P35609

- Molecule 2 is a protein called Myozenin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	45	374	243	64	64	3	0	0	0

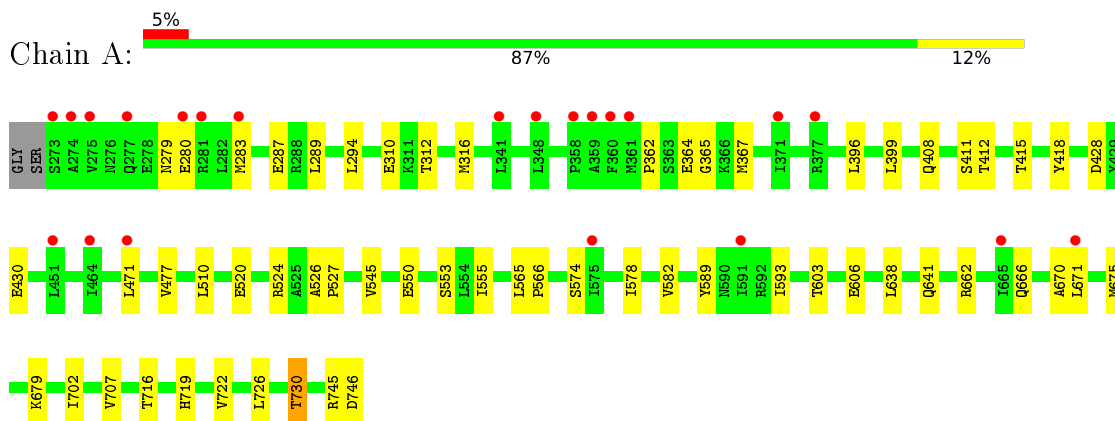
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	172	GLY	-	expression tag	UNP Q9NP98
B	173	PRO	-	expression tag	UNP Q9NP98
B	174	ALA	-	expression tag	UNP Q9NP98
B	175	MET	-	expression tag	UNP Q9NP98

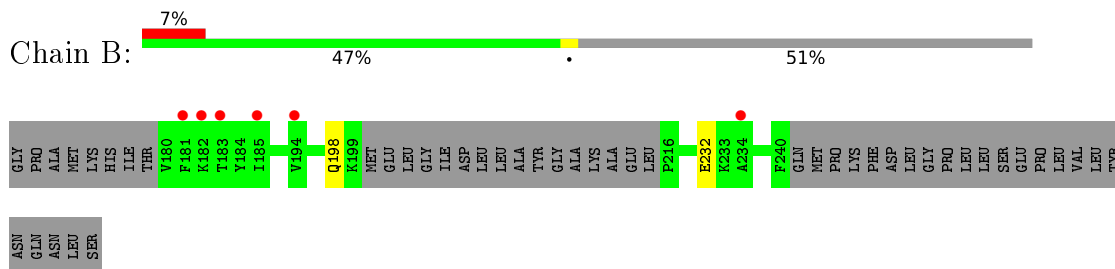
### 3 Residue-property plots [i](#)

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: Alpha-actinin-2



- Molecule 2: Myozenin-1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.44Å 103.44Å 218.14Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.58 – 2.69 46.73 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.58-2.69) 99.5 (46.73-2.69)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.88 (at 2.69Å)	Xtrriage
Refinement program	BUSTER 2.10.3	Depositor
R, $R_{free}$	0.204 , 0.235 0.223 , 0.258	Depositor DCC
$R_{free}$ test set	1862 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	80.5	Xtrriage
Anisotropy	0.717	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 87.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4309	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	125.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.61% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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	0.50	0/4001	0.65	0/5393
2	B	0.53	0/385	0.63	0/513
All	All	0.51	0/4386	0.64	0/5906

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	3935	0	3884	25	0
2	B	374	0	371	0	0
All	All	4309	0	4255	25	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 (25) 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:362:PRO:HG2	1:A:367:MET:HA	1.78	0.66
1:A:726:LEU:O	1:A:730:THR:HG23	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:GLU:H	1:A:316:MET:HE3	1.66	0.61
1:A:362:PRO:HD2	1:A:367:MET:HG2	1.93	0.50
1:A:716:THR:HG22	1:A:719:HIS:CG	2.47	0.50
1:A:510:LEU:HD13	1:A:589:TYR:HB3	1.96	0.48
1:A:280:GLU:HA	1:A:283:MET:HB2	1.95	0.48
1:A:415:THR:HA	1:A:418:TYR:CD2	2.50	0.47
1:A:603:THR:HG23	1:A:606:GLU:H	1.79	0.46
1:A:399:LEU:HD21	1:A:471:LEU:HB2	1.98	0.46
1:A:526:ALA:HB3	1:A:527:PRO:HD3	1.98	0.45
1:A:716:THR:HG22	1:A:719:HIS:ND1	2.32	0.45
1:A:671:LEU:HD23	1:A:745:ARG:HD3	1.98	0.45
1:A:675:MET:O	1:A:679:LYS:HB2	2.17	0.44
1:A:520:GLU:HG3	1:A:578:ILE:HD13	2.00	0.43
1:A:545:VAL:HG12	1:A:550:GLU:HB3	2.00	0.43
1:A:364:GLU:HB3	1:A:365:GLY:H	1.64	0.42
1:A:638:LEU:HA	1:A:641:GLN:HG2	2.02	0.42
1:A:316:MET:HG3	1:A:396:LEU:CD1	2.50	0.41
1:A:662:ARG:O	1:A:666:GLN:HB3	2.20	0.41
1:A:702:ILE:HG23	1:A:707:VAL:HB	2.03	0.41
1:A:408:GLN:O	1:A:412:THR:HG23	2.21	0.40
1:A:524:ARG:HH22	1:A:574:SER:HB3	1.86	0.40
1:A:565:LEU:N	1:A:566:PRO:HD2	2.37	0.40
1:A:411:SER:O	1:A:415:THR:HG23	2.22	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	472/476 (99%)	453 (96%)	18 (4%)	1 (0%)	47 73
2	B	41/92 (45%)	39 (95%)	2 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	513/568 (90%)	492 (96%)	20 (4%)	1 (0%)	47 73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	670	ALA

### 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	425/426 (100%)	410 (96%)	15 (4%)	36 65
2	B	39/78 (50%)	37 (95%)	2 (5%)	24 50
All	All	464/504 (92%)	447 (96%)	17 (4%)	34 63

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	279	ASN
1	A	287	GLU
1	A	289	LEU
1	A	294	LEU
1	A	312	THR
1	A	428	ASP
1	A	430	GLU
1	A	477	VAL
1	A	553	SER
1	A	555	ILE
1	A	582	VAL
1	A	593	ILE
1	A	722	VAL
1	A	730	THR
1	A	746	ASP
2	B	198	GLN
2	B	232	GLU



Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	317	GLN
1	A	426	GLN
1	A	455	GLN
1	A	480	ASN
1	A	616	GLN
1	A	641	GLN
1	A	680	GLN
1	A	684	ASN
1	A	712	HIS
2	B	222	ASN

### 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 monosaccharides 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

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	474/476 (99%)	0.37	22 (4%) 32 31	82, 115, 184, 226	0
2	B	45/92 (48%)	0.74	6 (13%) 3 2	100, 158, 192, 201	0
All	All	519/568 (91%)	0.41	28 (5%) 25 24	82, 117, 188, 226	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	181	PHE	5.7
1	A	359	ALA	5.5
1	A	591	ILE	5.4
1	A	358	PRO	4.3
1	A	277	GLN	4.2
1	A	665	ILE	4.0
1	A	281	ARG	4.0
2	B	183	THR	3.4
1	A	273	SER	3.4
1	A	280	GLU	3.1
1	A	341	LEU	3.1
1	A	274	ALA	3.1
1	A	377	ARG	2.7
1	A	671	LEU	2.5
1	A	275	VAL	2.5
2	B	194	VAL	2.4
1	A	283	MET	2.3
1	A	361	MET	2.3
1	A	464	ILE	2.3
2	B	182	LYS	2.3
1	A	471	LEU	2.2
1	A	348	LEU	2.2
1	A	575	ILE	2.2
1	A	451	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	185	ILE	2.1
1	A	360	PHE	2.1
2	B	234	ALA	2.1
1	A	371	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 monosaccharides 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.