



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 1, 2024 – 08:14 AM EDT

PDB ID : 8TF7
Title : Apo structure of protein crystal of Tri17
Authors : Zhai, R.; Zhang, W.
Deposited on : 2023-07-09
Resolution : 2.45 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.37.1
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) : 2.37.1

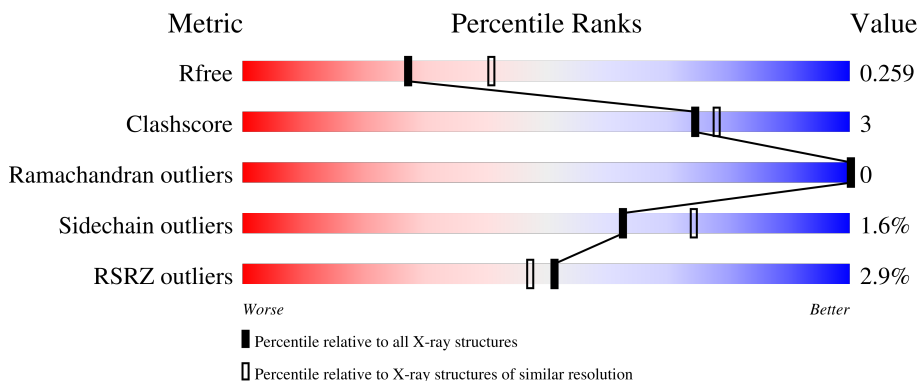
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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	565	 3% 83% 8% 9%
1	B	565	 3% 85% 7% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15997 atoms, of which 7758 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 AMP-binding protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	516	7859	2528	3878	699	741	13	0	0	0
1	B	517	7865	2530	3880	700	742	13	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A5H2UY12
A	557	LEU	-	expression tag	UNP A0A5H2UY12
A	558	GLU	-	expression tag	UNP A0A5H2UY12
A	559	HIS	-	expression tag	UNP A0A5H2UY12
A	560	HIS	-	expression tag	UNP A0A5H2UY12
A	561	HIS	-	expression tag	UNP A0A5H2UY12
A	562	HIS	-	expression tag	UNP A0A5H2UY12
A	563	HIS	-	expression tag	UNP A0A5H2UY12
A	564	HIS	-	expression tag	UNP A0A5H2UY12
A	565	END	-	expression tag	UNP A0A5H2UY12
B	1	MET	-	initiating methionine	UNP A0A5H2UY12
B	557	LEU	-	expression tag	UNP A0A5H2UY12
B	558	GLU	-	expression tag	UNP A0A5H2UY12
B	559	HIS	-	expression tag	UNP A0A5H2UY12
B	560	HIS	-	expression tag	UNP A0A5H2UY12
B	561	HIS	-	expression tag	UNP A0A5H2UY12
B	562	HIS	-	expression tag	UNP A0A5H2UY12
B	563	HIS	-	expression tag	UNP A0A5H2UY12
B	564	HIS	-	expression tag	UNP A0A5H2UY12
B	565	END	-	expression tag	UNP A0A5H2UY12

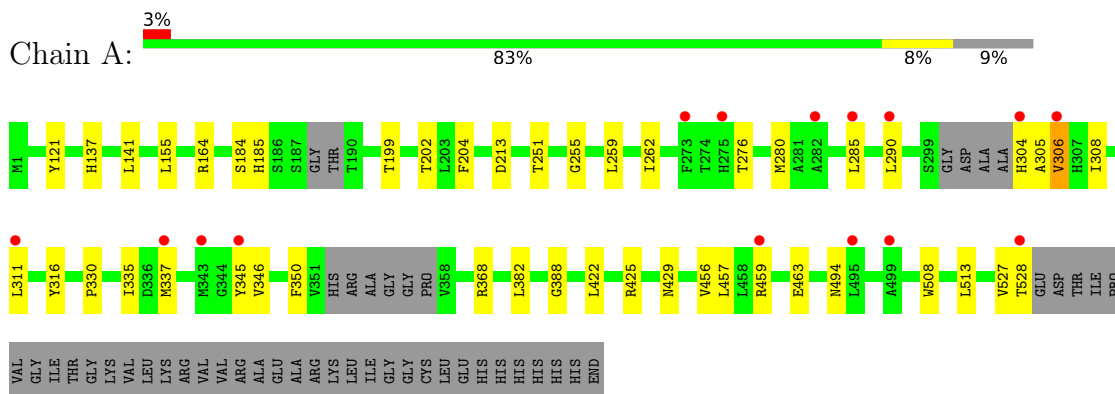
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	127	Total 127	O 127	0	0
2	B	146	Total 146	O 146	0	0

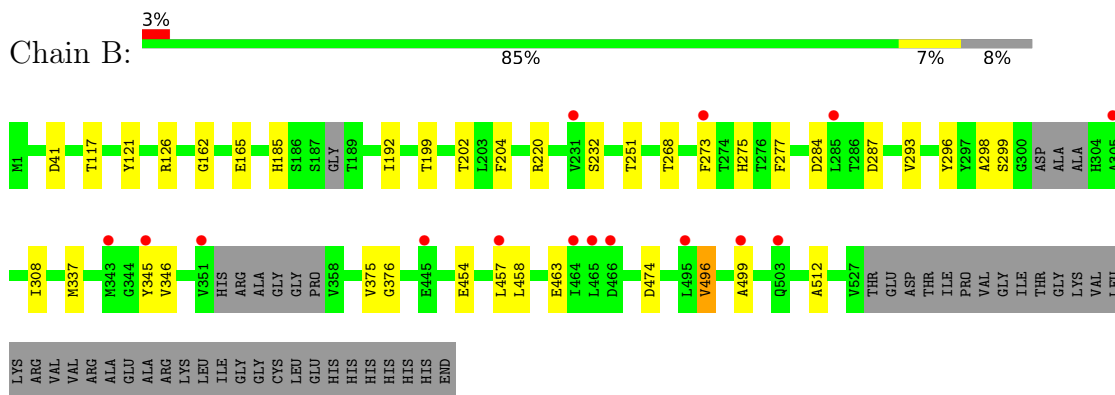
3 Residue-property plots [i](#)

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: AMP-binding protein



- Molecule 1: AMP-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	105.99Å 105.99Å 235.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.34 – 2.45 48.34 – 2.45	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.34-2.45) 100.0 (48.34-2.45)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.45Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.216 , 0.261 0.215 , 0.259	Depositor DCC
R_{free} test set	2525 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	39.6	Xtrriage
Anisotropy	0.443	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15997	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.06% 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	0.28	0/4077	0.52	0/5550
1	B	0.27	0/4081	0.51	0/5555
All	All	0.27	0/8158	0.52	0/11105

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	3981	3878	3905	27	0
1	B	3985	3880	3908	24	0
2	A	127	0	0	0	0
2	B	146	0	0	1	0
All	All	8239	7758	7813	51	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 (51) 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:496:VAL:HG23	1:B:499:ALA:HB2	1.42	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:345:TYR:CE2	1:B:346:VAL:HG22	2.13	0.83
1:B:345:TYR:CD2	1:B:346:VAL:HG22	2.18	0.77
1:B:463:GLU:O	1:B:496:VAL:HG22	1.91	0.70
1:A:494:ASN:OD1	1:A:527:VAL:CG2	2.41	0.68
1:B:277:PHE:HB3	1:B:308:ILE:HD11	1.79	0.65
1:A:305:ALA:HB1	1:A:308:ILE:HD11	1.79	0.62
1:A:285:LEU:HD13	1:A:311:LEU:HD23	1.87	0.57
1:A:494:ASN:OD1	1:A:527:VAL:HG21	2.05	0.55
1:A:137:HIS:CE1	1:A:141:LEU:HD11	2.45	0.52
1:A:255:GLY:O	1:A:259:LEU:HG	2.09	0.52
1:A:337:MET:SD	1:A:346:VAL:HG21	2.49	0.52
1:B:273:PHE:HB3	1:B:275:HIS:CD2	2.45	0.51
1:A:259:LEU:CD2	1:A:280:MET:HE3	2.41	0.51
1:B:375:VAL:HG12	1:B:376:GLY:O	2.11	0.49
1:B:162:GLY:O	1:B:165:GLU:HG2	2.13	0.49
1:A:457:LEU:HD21	1:A:513:LEU:HD21	1.96	0.47
1:A:316:TYR:CD1	1:A:330:PRO:HA	2.49	0.47
1:A:276:THR:O	1:A:280:MET:HG3	2.15	0.46
1:A:345:TYR:CG	1:A:346:VAL:N	2.84	0.45
1:A:382:LEU:O	1:A:425:ARG:NH1	2.48	0.45
1:A:527:VAL:HG23	1:A:528:THR:N	2.31	0.45
1:A:262:ILE:HG21	1:A:290:LEU:HD22	1.99	0.45
1:B:199:THR:OG1	1:B:202:THR:HG23	2.17	0.45
1:A:305:ALA:O	1:A:306:VAL:HG13	2.18	0.44
1:A:463:GLU:OE1	1:A:508:TRP:HZ2	2.01	0.43
1:B:273:PHE:HB2	1:B:275:HIS:NE2	2.33	0.43
1:A:305:ALA:HB1	1:A:308:ILE:CD1	2.45	0.43
1:A:121:TYR:CZ	1:A:185:HIS:HB3	2.53	0.43
1:B:277:PHE:HB3	1:B:308:ILE:CD1	2.47	0.43
1:A:388:GLY:O	1:A:422:LEU:HD23	2.18	0.43
1:A:456:VAL:HA	1:A:459:ARG:HD2	2.01	0.43
1:B:273:PHE:HB3	1:B:275:HIS:HD2	1.81	0.43
1:A:316:TYR:HD1	1:A:330:PRO:HA	1.82	0.43
1:A:199:THR:OG1	1:A:202:THR:HG23	2.19	0.43
1:B:126:ARG:NH1	2:B:605:HOH:O	2.53	0.42
1:B:220:ARG:HD2	1:B:268:THR:HG21	2.01	0.42
1:B:454:GLU:O	1:B:458:LEU:HD12	2.20	0.41
1:A:213:ASP:OD2	1:A:368:ARG:NE	2.51	0.41
1:A:137:HIS:CD2	1:A:155:LEU:HG	2.55	0.41
1:B:121:TYR:CZ	1:B:185:HIS:HB3	2.55	0.41
1:B:298:ALA:O	1:B:337:MET:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ILE:HG23	1:A:350:PHE:O	2.21	0.41
1:B:273:PHE:CB	1:B:275:HIS:CD2	3.03	0.41
1:B:293:VAL:HG11	1:B:296:TYR:CZ	2.56	0.41
1:B:345:TYR:CG	1:B:346:VAL:N	2.89	0.41
1:B:457:LEU:HD23	1:B:512:ALA:HB3	2.03	0.41
1:B:273:PHE:CE1	1:B:299:SER:HB3	2.56	0.40
1:A:388:GLY:C	1:A:422:LEU:HD23	2.41	0.40
1:B:117:THR:HG23	1:B:192:ILE:HD13	2.02	0.40
1:B:41:ASP:OD2	1:B:251:THR:HG22	2.21	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	508/565 (90%)	491 (97%)	17 (3%)	0	100	100
1	B	509/565 (90%)	493 (97%)	16 (3%)	0	100	100
All	All	1017/1130 (90%)	984 (97%)	33 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	412/447 (92%)	405 (98%)	7 (2%)	60	73
1	B	412/447 (92%)	406 (98%)	6 (2%)	65	76
All	All	824/894 (92%)	811 (98%)	13 (2%)	62	74

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	164	ARG
1	A	184	SER
1	A	204	PHE
1	A	251	THR
1	A	304	HIS
1	A	306	VAL
1	A	429	ASN
1	B	204	PHE
1	B	232	SER
1	B	284	ASP
1	B	287	ASP
1	B	474	ASP
1	B	496	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	307	HIS
1	B	275	HIS

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, 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	516/565 (91%)	0.06	15 (2%) 51 47	27, 47, 88, 123	0
1	B	517/565 (91%)	0.04	15 (2%) 51 47	26, 47, 85, 147	0
All	All	1033/1130 (91%)	0.05	30 (2%) 51 47	26, 47, 87, 147	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	345	TYR	9.8
1	B	345	TYR	7.2
1	A	306	VAL	4.5
1	A	459	ARG	4.0
1	A	282	ALA	4.0
1	B	495	LEU	3.5
1	B	285	LEU	3.4
1	A	285	LEU	3.4
1	A	495	LEU	3.3
1	A	273	PHE	3.3
1	A	275	HIS	3.2
1	A	528	THR	3.2
1	B	343	MET	3.1
1	B	273	PHE	3.0
1	B	445	GLU	3.0
1	A	499	ALA	2.9
1	A	304	HIS	2.7
1	B	464	ILE	2.6
1	B	351	VAL	2.5
1	B	499	ALA	2.5
1	A	311	LEU	2.4
1	B	465	LEU	2.4
1	B	457	LEU	2.3
1	B	305	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	231	VAL	2.3
1	A	337	MET	2.2
1	A	343	MET	2.2
1	B	503	GLN	2.1
1	A	290	LEU	2.1
1	B	466	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 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.