



wwPDB X-ray Structure Validation Summary Report ⓘ

May 17, 2020 – 05:40 am BST

PDB ID : 1TIM
Title : STRUCTURE OF TRIOSE PHOSPHATE ISOMERASE FROM CHICKEN MUSCLE
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Deposited on : 1976-09-01
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

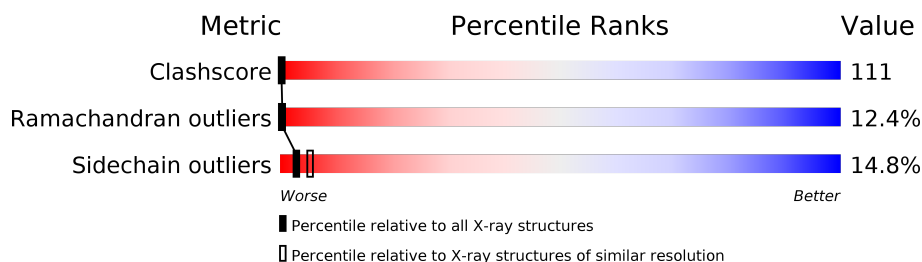
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

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 on 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	247	
1	B	247	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3740 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 TRIOSEPHOSPHATE ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	0	0
			1870	1187	329	348	6			
1	B	247	Total	C	N	O	S	0	0	0
			1870	1187	329	348	6			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	LYS	ASP	CONFLICT	UNP P00940
A	18	ARG	LYS	CONFLICT	UNP P00940
A	29	ASP	ASN	CONFLICT	UNP P00940
A	145	GLN	GLU	CONFLICT	UNP P00940
A	146	GLU	GLN	CONFLICT	UNP P00940
A	194	THR	SER	CONFLICT	UNP P00940
A	202	VAL	-	INSERTION	UNP P00940
A	?	-	THR	DELETION	UNP P00940
B	17	LYS	ASP	CONFLICT	UNP P00940
B	18	ARG	LYS	CONFLICT	UNP P00940
B	29	ASP	ASN	CONFLICT	UNP P00940
B	145	GLN	GLU	CONFLICT	UNP P00940
B	146	GLU	GLN	CONFLICT	UNP P00940
B	194	THR	SER	CONFLICT	UNP P00940
B	202	VAL	-	INSERTION	UNP P00940
B	?	-	THR	DELETION	UNP P00940

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	106.01Å 74.76Å 61.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	unknown	Depositor
R, R_{free}	(Not available) , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3740	wwPDB-VP
Average B, all atoms (Å ²)	0.0	wwPDB-VP

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	1.03	9/1906 (0.5%)	1.53	14/2574 (0.5%)
1	B	1.06	9/1906 (0.5%)	1.70	32/2574 (1.2%)
All	All	1.05	18/3812 (0.5%)	1.62	46/5148 (0.9%)

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	134	ARG	NE-CZ	12.41	1.49	1.33
1	A	189	ARG	NE-CZ	11.05	1.47	1.33
1	A	52	ARG	NE-CZ	11.01	1.47	1.33
1	A	18	ARG	NE-CZ	10.99	1.47	1.33
1	B	99	ARG	NE-CZ	10.95	1.47	1.33

The worst 5 of 46 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	134	ARG	NE-CZ-NH1	13.62	127.11	120.30
1	B	156	ASP	CB-CG-OD2	-13.20	106.42	118.30
1	B	156	ASP	CB-CG-OD1	12.42	129.48	118.30
1	B	134	ARG	CD-NE-CZ	-10.27	109.23	123.60
1	B	217	CYS	N-CA-CB	-9.93	92.73	110.60

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	1870	0	1887	474	2
1	B	1870	0	1886	420	2
All	All	3740	0	3773	834	2

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 111.

The worst 5 of 834 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:45:SER:C	1:B:46:ILE:HG22	1.44	1.38
1:A:46:ILE:CG2	1:B:45:SER:O	1.76	1.32
1:A:45:SER:O	1:A:48:LEU:HB2	1.16	1.28
1:B:129:GLU:OE1	1:B:134:ARG:HB2	1.32	1.27
1:A:189:ARG:NE	1:A:225:ASP:OD2	1.67	1.27

All (2) 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:180:GLN:NE2	1:B:171:GLY:O[2_565]	1.08	1.12
1:A:180:GLN:NE2	1:B:171:GLY:C[2_565]	1.90	0.30

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	245/247 (99%)	144 (59%)	63 (26%)	38 (16%)	0	0
1	B	245/247 (99%)	162 (66%)	60 (24%)	23 (9%)	0	0
All	All	490/494 (99%)	306 (62%)	123 (25%)	61 (12%)	0	0

5 of 61 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	31	ALA
1	A	57	ALA
1	A	58	LYS
1	A	77	GLU
1	A	104	GLU

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	192/192 (100%)	163 (85%)	29 (15%)	3	5
1	B	192/192 (100%)	164 (85%)	28 (15%)	3	5
All	All	384/384 (100%)	327 (85%)	57 (15%)	3	5

5 of 57 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	229	PHE
1	B	24	LEU
1	B	229	PHE
1	A	231	VAL
1	A	242	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	203	GLN
1	B	245	ASN
1	A	245	ASN
1	A	145	GLN
1	A	223	GLN

5.3.3 RNA ⓘ

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 carbohydrates 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	98:ARG	C	99:ARG	N	1.19

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.