



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 11:17 am BST

PDB ID : 1YJ8
Title : Initial structural analysis of Plasmodium falciparum Glycerol-3-phosphate dehydrogenase
Authors : Robien, M.A.; Hol, W.G.J.; Structural Genomics of Pathogenic Protozoa Consortium (SGPP)
Deposited on : 2005-01-13
Resolution : 2.85 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 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	:	2.11
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)	:	2.11

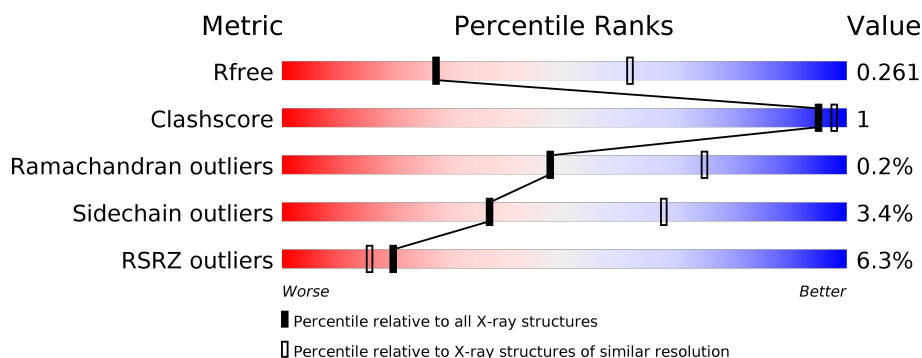
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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 $\leq 5\%$. 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	375	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>5%</div> </div> </div>
1	B	375	<div> <div>8%</div> <div> <div></div> <div>87%</div> <div>8%</div> <div>5%</div> </div> </div>
1	C	375	<div> <div>8%</div> <div> <div></div> <div>89%</div> <div>6%</div> <div>5%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8425 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 glycerol-3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	0	0	0
			2812	1806	466	522	18			
1	B	357	Total	C	N	O	S	0	0	0
			2806	1802	464	522	18			
1	C	357	Total	C	N	O	S	0	0	0
			2797	1796	462	521	18			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	CLONING ARTIFACT	UNP Q8I5P5
A	2	ALA	-	CLONING ARTIFACT	UNP Q8I5P5
A	3	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
A	4	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
A	5	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
A	6	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
A	7	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
A	8	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
B	1	MET	-	CLONING ARTIFACT	UNP Q8I5P5
B	2	ALA	-	CLONING ARTIFACT	UNP Q8I5P5
B	3	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
B	4	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
B	5	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
B	6	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
B	7	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
B	8	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
C	1	MET	-	CLONING ARTIFACT	UNP Q8I5P5
C	2	ALA	-	CLONING ARTIFACT	UNP Q8I5P5
C	3	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
C	4	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
C	5	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
C	6	HIS	-	CLONING ARTIFACT	UNP Q8I5P5
C	7	HIS	-	CLONING ARTIFACT	UNP Q8I5P5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	8	HIS	-	CLONING ARTIFACT	UNP Q8I5P5

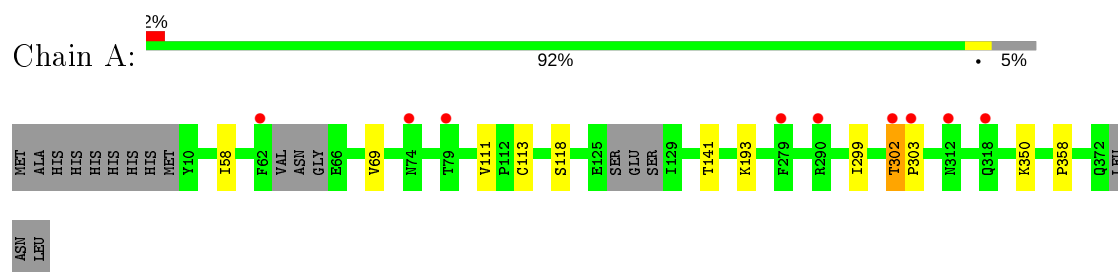
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total O 4 4	0	0
2	B	3	Total O 3 3	0	0
2	C	3	Total O 3 3	0	0

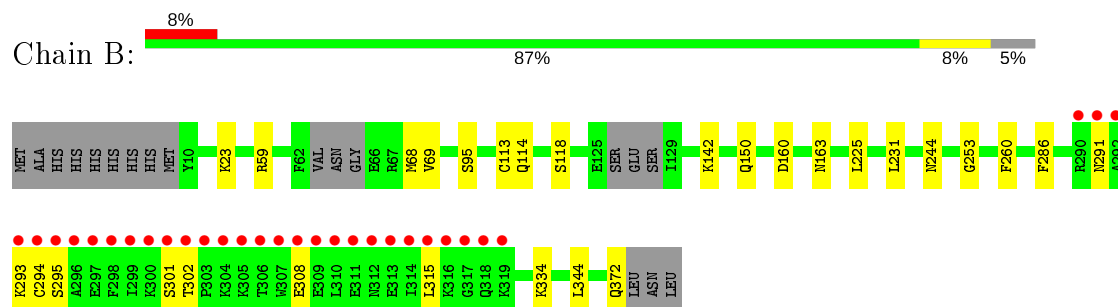
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

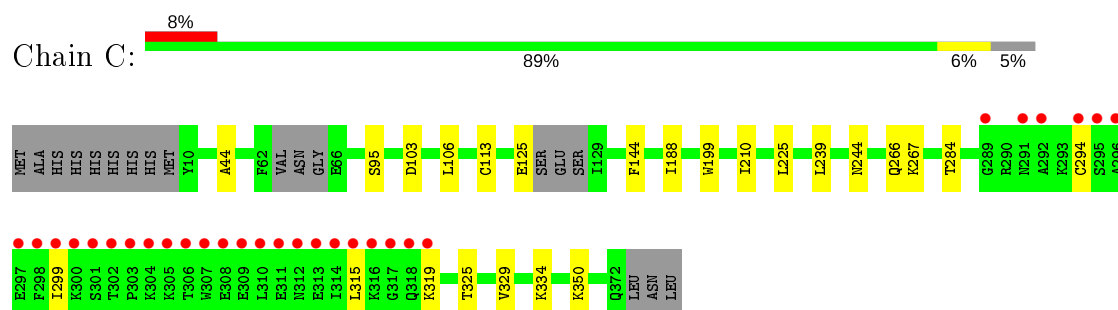
- Molecule 1: glycerol-3-phosphate dehydrogenase



- Molecule 1: glycerol-3-phosphate dehydrogenase



- Molecule 1: glycerol-3-phosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	177.46 Å 177.46 Å 232.70 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.54 – 2.85 44.54 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.0 (44.54-2.85) 99.1 (44.54-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.86 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.236 , 0.256 0.240 , 0.261	Depositor DCC
R_{free} test set	2557 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	58.7	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.3	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	0.91	EDS
Total number of atoms	8425	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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	0.36	0/2862	0.50	0/3861
1	B	0.37	0/2856	0.50	0/3855
1	C	0.37	0/2847	0.49	0/3844
All	All	0.37	0/8565	0.50	0/11560

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	2812	0	2844	7	0
1	B	2806	0	2826	6	0
1	C	2797	0	2808	8	0
2	A	4	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	1	0
All	All	8425	0	8478	21	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 1.

All (21) 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:111:VAL:O	1:A:141:THR:HG21	2.02	0.59
1:A:302:THR:HB	1:A:303:PRO:HD3	1.85	0.59
1:A:58:ILE:HD11	1:A:69:VAL:HA	1.87	0.56
1:C:144:PHE:HE1	1:C:325:THR:HG22	1.73	0.53
1:C:294:CYS:SG	1:C:315:LEU:HG	2.49	0.52
1:A:302:THR:CB	1:A:303:PRO:HD3	2.41	0.51
1:B:69:VAL:HG21	1:B:95:SER:HB3	1.95	0.48
1:C:44:ALA:O	2:C:501:HOH:O	2.20	0.47
1:C:239:LEU:HG	1:C:299:ILE:HG21	1.97	0.45
1:A:302:THR:CB	1:A:303:PRO:CD	2.95	0.45
1:B:294:CYS:SG	1:B:315:LEU:HD21	2.57	0.45
1:B:253:GLY:HA2	1:B:344:LEU:HD23	1.99	0.44
1:C:188:ILE:HD12	1:C:210:ILE:HG21	2.00	0.43
1:B:225:LEU:HD11	1:B:260:PHE:CD2	2.53	0.43
1:A:111:VAL:C	1:A:141:THR:HG21	2.40	0.42
1:C:225:LEU:HD22	1:C:329:VAL:HG21	2.02	0.41
1:B:231:LEU:HD12	1:B:295:SER:HB3	2.02	0.41
1:B:68:MET:HE2	1:B:68:MET:HB2	1.98	0.41
1:C:144:PHE:CE1	1:C:325:THR:HG22	2.53	0.40
1:A:299:ILE:HD11	1:A:358:PRO:HB3	2.03	0.40
1:C:106:LEU:HB3	1:C:199:TRP:CH2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	351/375 (94%)	342 (97%)	8 (2%)	1 (0%)	41	68
1	B	351/375 (94%)	338 (96%)	12 (3%)	1 (0%)	41	68
1	C	351/375 (94%)	341 (97%)	10 (3%)	0	100	100
All	All	1053/1125 (94%)	1021 (97%)	30 (3%)	2 (0%)	47	75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	302	THR
1	B	291	ASN

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	314/333 (94%)	310 (99%)	4 (1%)	69	88
1	B	312/333 (94%)	295 (95%)	17 (5%)	22	49
1	C	310/333 (93%)	299 (96%)	11 (4%)	36	67
All	All	936/999 (94%)	904 (97%)	32 (3%)	37	67

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	CYS
1	A	118	SER
1	A	193	LYS
1	A	350	LYS
1	B	23	LYS
1	B	59	ARG
1	B	113	CYS
1	B	114	GLN
1	B	118	SER
1	B	142	LYS
1	B	150	GLN
1	B	160	ASP
1	B	163	ASN
1	B	244	ASN
1	B	286	PHE
1	B	293	LYS
1	B	301	SER
1	B	302	THR
1	B	308	GLU

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Mol	Chain	Res	Type
1	B	334	LYS
1	B	372	GLN
1	C	95	SER
1	C	103	ASP
1	C	113	CYS
1	C	125	GLU
1	C	244	ASN
1	C	266	GLN
1	C	267	LYS
1	C	284	THR
1	C	319	LYS
1	C	334	LYS
1	C	350	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 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 ⓘ

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	357/375 (95%)	0.11	9 (2%) 57 54	41, 55, 78, 92	2 (0%)
1	B	357/375 (95%)	0.28	30 (8%) 11 7	39, 56, 105, 113	0
1	C	357/375 (95%)	0.39	29 (8%) 12 8	39, 56, 108, 115	0
All	All	1071/1125 (95%)	0.26	68 (6%) 20 15	39, 55, 91, 115	2 (0%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	302	THR	9.7
1	C	307	TRP	9.3
1	B	303	PRO	8.8
1	C	303	PRO	8.5
1	C	310	LEU	8.1
1	B	302	THR	7.3
1	C	301	SER	6.9
1	B	299	ILE	6.6
1	C	318	GLN	6.2
1	B	314	ILE	5.8
1	C	306	THR	5.6
1	B	310	LEU	5.5
1	B	292	ALA	5.4
1	C	304	LYS	5.4
1	C	314	ILE	5.3
1	B	300	LYS	5.3
1	C	312	ASN	5.2
1	C	309	GLU	5.1
1	B	294	CYS	5.1
1	B	318	GLN	5.0
1	B	309	GLU	4.9
1	C	311	GLU	4.9
1	C	299	ILE	4.9

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Mol	Chain	Res	Type	RSRZ
1	B	317	GLY	4.8
1	B	301	SER	4.7
1	C	305	LYS	4.7
1	B	307	TRP	4.7
1	B	304	LYS	4.6
1	B	305	LYS	4.3
1	C	300	LYS	4.2
1	C	315	LEU	4.1
1	C	292	ALA	4.0
1	B	306	THR	4.0
1	C	308	GLU	4.0
1	B	312	ASN	3.7
1	B	319	LYS	3.6
1	C	296	ALA	3.5
1	B	313	GLU	3.5
1	B	308	GLU	3.5
1	B	293	LYS	3.5
1	C	294	CYS	3.4
1	C	316	LYS	3.4
1	C	317	GLY	3.4
1	C	319	LYS	3.3
1	A	302	THR	3.3
1	B	297	GLU	3.3
1	B	290	ARG	3.3
1	C	298	PHE	3.2
1	B	316	LYS	3.1
1	B	315	LEU	2.9
1	B	296	ALA	2.8
1	C	291	ASN	2.6
1	B	291	ASN	2.6
1	C	289	GLY	2.6
1	C	313	GLU	2.5
1	A	279	PHE	2.4
1	C	297	GLU	2.4
1	B	295	SER	2.2
1	A	79	THR	2.2
1	A	318	GLN	2.2
1	A	74	ASN	2.1
1	C	295	SER	2.1
1	A	303	PRO	2.1
1	A	290	ARG	2.1
1	B	298	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	311	GLU	2.0
1	A	62	PHE	2.0
1	A	312	ASN	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 carbohydrates 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.