



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 05:52 am BST

PDB ID : 3IG0
Title : crystal structure of the second part of the Mycobacterium tuberculosis DNA gyrase reaction core: the TOPRIM domain at 2.1 Å resolution
Authors : Piton, J.; Aubry, A.; Delarue, M.; Mayer, C.
Deposited on : 2009-07-27
Resolution : 2.10 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

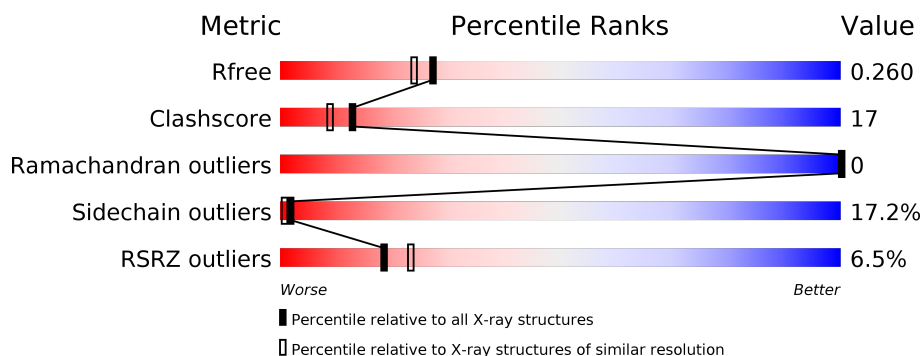
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	242	<div> <div>5%</div> <div>52%</div> <div>19%</div> <div>5%</div> <div>24%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1581 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 DNA gyrase subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	185	Total	C	N	O	S	0	0	0
			1474	940	253	275	6			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	434	MET	-	EXPRESSION TAG	UNP P0C5C5
A	435	ALA	-	EXPRESSION TAG	UNP P0C5C5
A	436	HIS	-	EXPRESSION TAG	UNP P0C5C5
A	437	HIS	-	EXPRESSION TAG	UNP P0C5C5
A	438	HIS	-	EXPRESSION TAG	UNP P0C5C5
A	439	HIS	-	EXPRESSION TAG	UNP P0C5C5
A	440	HIS	-	EXPRESSION TAG	UNP P0C5C5
A	441	HIS	-	EXPRESSION TAG	UNP P0C5C5
A	442	VAL	-	EXPRESSION TAG	UNP P0C5C5
A	443	ASP	-	EXPRESSION TAG	UNP P0C5C5
A	444	ASP	-	EXPRESSION TAG	UNP P0C5C5
A	445	ASP	-	EXPRESSION TAG	UNP P0C5C5
A	446	ASP	-	EXPRESSION TAG	UNP P0C5C5
A	447	LYS	-	EXPRESSION TAG	UNP P0C5C5

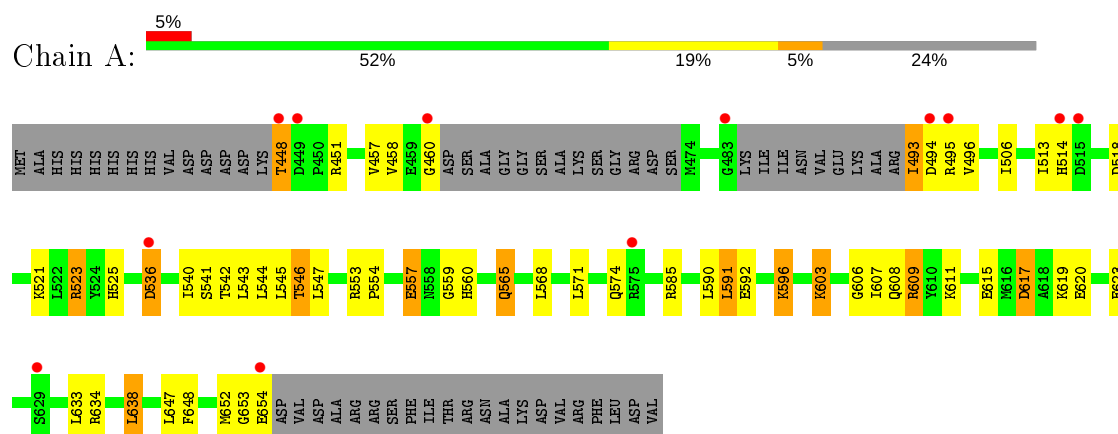
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	107	Total	O	0	0
			107	107		

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: DNA gyrase subunit B



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	52.87Å 52.87Å 190.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	13.96 – 2.10 13.96 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.6 (13.96-2.10) 99.6 (13.96-2.10)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.10Å)	Xtriage
Refinement program	BUSTER-TNT 2.5.1	Depositor
R, R_{free}	0.214 , 0.249 0.229 , 0.260	Depositor DCC
R_{free} test set	799 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	33.5	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 61.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	1581	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.44% 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

5.1 Standard geometry

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	0.32	0/1498	0.53	1/2017 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	460	GLY	N-CA-C	-5.34	99.74	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1474	0	1492	49	0
2	A	107	0	0	3	0
All	All	1581	0	1492	49	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 17.

All (49) 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:565:GLN:HG2	1:A:634:ARG:HH22	1.20	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:565:GLN:HG3	1:A:634:ARG:HH12	1.27	0.99
1:A:609:ARG:HH11	1:A:609:ARG:HG2	1.43	0.84
1:A:619:LYS:O	1:A:623:GLU:HG3	1.84	0.78
1:A:603:LYS:HZ2	1:A:603:LYS:HB3	1.53	0.72
1:A:565:GLN:HG2	1:A:634:ARG:NH2	2.02	0.72
1:A:523:ARG:HG2	2:A:130:HOH:O	1.90	0.71
1:A:565:GLN:HG3	1:A:634:ARG:NH1	2.05	0.71
1:A:496:VAL:HG11	1:A:543:LEU:HD11	1.75	0.68
1:A:536:ASP:OD1	1:A:536:ASP:N	2.28	0.66
1:A:607:ILE:N	1:A:607:ILE:HD12	2.11	0.65
1:A:542:THR:O	1:A:546:THR:HG23	1.96	0.65
1:A:565:GLN:CG	1:A:634:ARG:HH12	2.04	0.63
1:A:565:GLN:CG	1:A:634:ARG:HH22	2.04	0.63
1:A:596:LYS:HD3	1:A:596:LYS:N	2.13	0.62
1:A:617:ASP:HB3	1:A:620:GLU:HG3	1.82	0.61
1:A:523:ARG:HG3	2:A:121:HOH:O	2.00	0.60
1:A:653:GLY:O	1:A:654:GLU:HB2	2.02	0.60
1:A:617:ASP:OD1	1:A:620:GLU:HG3	2.01	0.60
1:A:546:THR:HG21	1:A:648:PHE:HB2	1.87	0.57
1:A:554:PRO:HA	1:A:557:GLU:HG3	1.87	0.56
1:A:603:LYS:HB3	1:A:603:LYS:NZ	2.19	0.56
1:A:553:ARG:O	1:A:557:GLU:HG3	2.06	0.55
1:A:609:ARG:NH1	1:A:609:ARG:HG2	2.16	0.53
1:A:457:VAL:HG11	1:A:544:LEU:CD2	2.40	0.52
1:A:606:GLY:C	1:A:607:ILE:HD12	2.31	0.51
1:A:546:THR:HG21	1:A:648:PHE:CB	2.41	0.51
1:A:609:ARG:CG	1:A:609:ARG:HH11	2.18	0.49
1:A:457:VAL:HG11	1:A:544:LEU:HD21	1.94	0.49
1:A:525:HIS:O	1:A:560:HIS:HD2	1.97	0.48
1:A:652:MET:C	1:A:654:GLU:N	2.66	0.48
1:A:493:ILE:HD13	1:A:493:ILE:HA	1.74	0.46
1:A:603:LYS:O	1:A:603:LYS:HG3	2.15	0.46
1:A:607:ILE:CD1	1:A:607:ILE:N	2.79	0.46
1:A:513:ILE:HG23	1:A:514:HIS:N	2.32	0.45
1:A:647:LEU:O	1:A:647:LEU:HD12	2.16	0.45
1:A:448:THR:OG1	1:A:448:THR:O	2.33	0.45
1:A:559:GLY:HA2	2:A:4:HOH:O	2.16	0.45
1:A:596:LYS:CD	1:A:596:LYS:N	2.80	0.44
1:A:518:ASP:HB3	1:A:521:LYS:HD2	1.99	0.44
1:A:648:PHE:O	1:A:652:MET:HG2	2.18	0.44
1:A:448:THR:O	1:A:523:ARG:NH2	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:652:MET:C	1:A:654:GLU:H	2.22	0.43
1:A:506:ILE:CG2	1:A:513:ILE:HD13	2.49	0.43
1:A:609:ARG:CG	1:A:609:ARG:NH1	2.77	0.43
1:A:545:LEU:HA	1:A:545:LEU:HD23	1.86	0.42
1:A:571:LEU:CD2	1:A:591:LEU:HD13	2.51	0.41
1:A:608:GLN:NE2	1:A:615:GLU:OE2	2.40	0.40
1:A:638:LEU:HD23	1:A:638:LEU:HA	1.80	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	179/242 (74%)	174 (97%)	5 (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	157/205 (77%)	130 (83%)	27 (17%)	2	1

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	448	THR
1	A	451	ARG
1	A	458	VAL
1	A	493	ILE
1	A	494	ASP
1	A	495	ARG
1	A	523	ARG
1	A	536	ASP
1	A	540	ILE
1	A	541	SER
1	A	546	THR
1	A	547	LEU
1	A	557	GLU
1	A	565	GLN
1	A	568	LEU
1	A	574	GLN
1	A	585	ARG
1	A	590	LEU
1	A	591	LEU
1	A	592	GLU
1	A	596	LYS
1	A	603	LYS
1	A	609	ARG
1	A	611	LYS
1	A	617	ASP
1	A	633	LEU
1	A	638	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	514	HIS
1	A	560	HIS
1	A	574	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](#)

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	185/242 (76%)	0.27	12 (6%) 18 23	18, 36, 59, 88	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	495	ARG	6.5
1	A	448	THR	6.4
1	A	575	ARG	4.8
1	A	654	GLU	3.6
1	A	460	GLY	3.0
1	A	494	ASP	2.9
1	A	483	GLY	2.8
1	A	629	SER	2.8
1	A	514	HIS	2.4
1	A	515	ASP	2.3
1	A	449	ASP	2.0
1	A	536	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 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.