



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 4, 2024 – 12:09 PM EDT

PDB ID : 8TXA
Title : Apo Structure of (N1G37) tRNA Methyltransferase from *Mycobacterium marinum*
Authors : Balsamo, A.; Bruno, C.; Edele, C.; Fabian, T.; Jannotta, R.; Lee, R.; Schryver, D.; Warsaw, J.; Warsaw, L.; Stojanoff, V.; Battaile, K.; Perez, A.; Bolen, R.
Deposited on : 2023-08-23
Resolution : 1.59 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 : 2.36.1
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) : 2.36.1

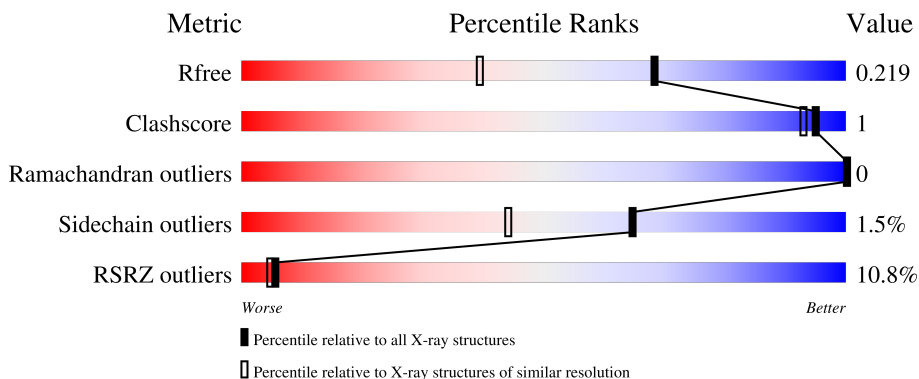
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.59 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	250	 5% 73% 24%
1	B	250	 8% 81% 17%
1	C	250	 8% 72% 25%
1	D	250	 12% 76% 19%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6420 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 tRNA (guanine-N(1)-)-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	190	Total 1445	C 915	N 258	O 265	S 7	0	0	0
1	B	207	Total 1570	C 989	N 284	O 290	S 7	0	0	0
1	C	187	Total 1417	C 898	N 250	O 262	S 7	0	0	0
1	D	203	Total 1529	C 962	N 275	O 285	S 7	0	1	0

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP A0A2Z5YCU9
A	-19	ALA	-	expression tag	UNP A0A2Z5YCU9
A	-18	HIS	-	expression tag	UNP A0A2Z5YCU9
A	-17	HIS	-	expression tag	UNP A0A2Z5YCU9
A	-16	HIS	-	expression tag	UNP A0A2Z5YCU9
A	-15	HIS	-	expression tag	UNP A0A2Z5YCU9
A	-14	HIS	-	expression tag	UNP A0A2Z5YCU9
A	-13	HIS	-	expression tag	UNP A0A2Z5YCU9
A	-12	MET	-	expression tag	UNP A0A2Z5YCU9
A	-11	GLY	-	expression tag	UNP A0A2Z5YCU9
A	-10	THR	-	expression tag	UNP A0A2Z5YCU9
A	-9	LEU	-	expression tag	UNP A0A2Z5YCU9
A	-8	GLU	-	expression tag	UNP A0A2Z5YCU9
A	-7	ALA	-	expression tag	UNP A0A2Z5YCU9
A	-6	GLN	-	expression tag	UNP A0A2Z5YCU9
A	-5	THR	-	expression tag	UNP A0A2Z5YCU9
A	-4	GLN	-	expression tag	UNP A0A2Z5YCU9
A	-3	GLY	-	expression tag	UNP A0A2Z5YCU9
A	-2	PRO	-	expression tag	UNP A0A2Z5YCU9
A	-1	GLY	-	expression tag	UNP A0A2Z5YCU9
A	0	SER	-	expression tag	UNP A0A2Z5YCU9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	MET	-	initiating methionine	UNP A0A2Z5YCU9
B	-19	ALA	-	expression tag	UNP A0A2Z5YCU9
B	-18	HIS	-	expression tag	UNP A0A2Z5YCU9
B	-17	HIS	-	expression tag	UNP A0A2Z5YCU9
B	-16	HIS	-	expression tag	UNP A0A2Z5YCU9
B	-15	HIS	-	expression tag	UNP A0A2Z5YCU9
B	-14	HIS	-	expression tag	UNP A0A2Z5YCU9
B	-13	HIS	-	expression tag	UNP A0A2Z5YCU9
B	-12	MET	-	expression tag	UNP A0A2Z5YCU9
B	-11	GLY	-	expression tag	UNP A0A2Z5YCU9
B	-10	THR	-	expression tag	UNP A0A2Z5YCU9
B	-9	LEU	-	expression tag	UNP A0A2Z5YCU9
B	-8	GLU	-	expression tag	UNP A0A2Z5YCU9
B	-7	ALA	-	expression tag	UNP A0A2Z5YCU9
B	-6	GLN	-	expression tag	UNP A0A2Z5YCU9
B	-5	THR	-	expression tag	UNP A0A2Z5YCU9
B	-4	GLN	-	expression tag	UNP A0A2Z5YCU9
B	-3	GLY	-	expression tag	UNP A0A2Z5YCU9
B	-2	PRO	-	expression tag	UNP A0A2Z5YCU9
B	-1	GLY	-	expression tag	UNP A0A2Z5YCU9
B	0	SER	-	expression tag	UNP A0A2Z5YCU9
C	-20	MET	-	initiating methionine	UNP A0A2Z5YCU9
C	-19	ALA	-	expression tag	UNP A0A2Z5YCU9
C	-18	HIS	-	expression tag	UNP A0A2Z5YCU9
C	-17	HIS	-	expression tag	UNP A0A2Z5YCU9
C	-16	HIS	-	expression tag	UNP A0A2Z5YCU9
C	-15	HIS	-	expression tag	UNP A0A2Z5YCU9
C	-14	HIS	-	expression tag	UNP A0A2Z5YCU9
C	-13	HIS	-	expression tag	UNP A0A2Z5YCU9
C	-12	MET	-	expression tag	UNP A0A2Z5YCU9
C	-11	GLY	-	expression tag	UNP A0A2Z5YCU9
C	-10	THR	-	expression tag	UNP A0A2Z5YCU9
C	-9	LEU	-	expression tag	UNP A0A2Z5YCU9
C	-8	GLU	-	expression tag	UNP A0A2Z5YCU9
C	-7	ALA	-	expression tag	UNP A0A2Z5YCU9
C	-6	GLN	-	expression tag	UNP A0A2Z5YCU9
C	-5	THR	-	expression tag	UNP A0A2Z5YCU9
C	-4	GLN	-	expression tag	UNP A0A2Z5YCU9
C	-3	GLY	-	expression tag	UNP A0A2Z5YCU9
C	-2	PRO	-	expression tag	UNP A0A2Z5YCU9
C	-1	GLY	-	expression tag	UNP A0A2Z5YCU9
C	0	SER	-	expression tag	UNP A0A2Z5YCU9

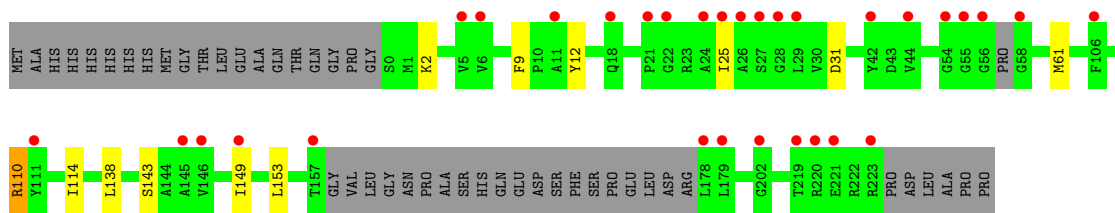
Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	-20	MET	-	initiating methionine	UNP A0A2Z5YCU9
D	-19	ALA	-	expression tag	UNP A0A2Z5YCU9
D	-18	HIS	-	expression tag	UNP A0A2Z5YCU9
D	-17	HIS	-	expression tag	UNP A0A2Z5YCU9
D	-16	HIS	-	expression tag	UNP A0A2Z5YCU9
D	-15	HIS	-	expression tag	UNP A0A2Z5YCU9
D	-14	HIS	-	expression tag	UNP A0A2Z5YCU9
D	-13	HIS	-	expression tag	UNP A0A2Z5YCU9
D	-12	MET	-	expression tag	UNP A0A2Z5YCU9
D	-11	GLY	-	expression tag	UNP A0A2Z5YCU9
D	-10	THR	-	expression tag	UNP A0A2Z5YCU9
D	-9	LEU	-	expression tag	UNP A0A2Z5YCU9
D	-8	GLU	-	expression tag	UNP A0A2Z5YCU9
D	-7	ALA	-	expression tag	UNP A0A2Z5YCU9
D	-6	GLN	-	expression tag	UNP A0A2Z5YCU9
D	-5	THR	-	expression tag	UNP A0A2Z5YCU9
D	-4	GLN	-	expression tag	UNP A0A2Z5YCU9
D	-3	GLY	-	expression tag	UNP A0A2Z5YCU9
D	-2	PRO	-	expression tag	UNP A0A2Z5YCU9
D	-1	GLY	-	expression tag	UNP A0A2Z5YCU9
D	0	SER	-	expression tag	UNP A0A2Z5YCU9

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	99	Total O 99 99	0	0
2	B	137	Total O 137 137	0	0
2	C	113	Total O 113 113	0	0
2	D	110	Total O 110 110	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.94Å 60.80Å 68.07Å 97.05° 111.99° 100.69°	Depositor
Resolution (Å)	22.98 – 1.59 22.98 – 1.59	Depositor EDS
% Data completeness (in resolution range)	59.5 (22.98-1.59) 59.4 (22.98-1.59)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 1.59Å)	Xtrriage
Refinement program	BUSTER 2.10.4	Depositor
R, R_{free}	0.188 , 0.224 0.183 , 0.219	Depositor DCC
R_{free} test set	3210 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	20.7	Xtrriage
Anisotropy	0.005	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6420	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.26% 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	0.50	0/1478	0.58	0/2016
1	B	0.53	0/1603	0.61	0/2183
1	C	0.54	0/1447	0.59	0/1972
1	D	0.51	0/1559	0.59	0/2123
All	All	0.52	0/6087	0.59	0/8294

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	1445	0	1402	3	0
1	B	1570	0	1534	2	0
1	C	1417	0	1367	4	0
1	D	1529	0	1480	7	0
2	A	99	0	0	0	0
2	B	137	0	0	0	0
2	C	113	0	0	0	0
2	D	110	0	0	0	0
All	All	6420	0	5783	14	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 (14) 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:D:9:PHE:HE2	1:D:110:ARG:HD3	1.61	0.65
1:D:12:TYR:OH	1:D:143:SER:HB3	2.01	0.60
1:C:198:ILE:HD11	1:C:206:ARG:HB3	1.84	0.58
1:B:12:TYR:OH	1:B:143:SER:HB3	2.05	0.56
1:D:9:PHE:CE2	1:D:110:ARG:HD3	2.41	0.53
1:C:30:VAL:HG21	1:C:156:LEU:HD23	1.90	0.52
1:C:198:ILE:CD1	1:C:206:ARG:HB3	2.39	0.52
1:A:62:LYS:HB3	1:A:64:PRO:HD2	1.91	0.52
1:A:2:LYS:HD2	1:A:31:ASP:O	2.12	0.49
1:D:2:LYS:HD2	1:D:31:ASP:O	2.16	0.46
1:A:50:ASP:OD2	1:B:186:ARG:HG2	2.16	0.45
1:D:61:MET:HB2	1:D:114:ILE:HA	2.01	0.43
1:C:147:VAL:HG13	1:D:138:LEU:HD21	2.02	0.41
1:D:149:ILE:O	1:D:153:LEU:HG	2.21	0.41

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	184/250 (74%)	183 (100%)	1 (0%)	0	100	100
1	B	201/250 (80%)	200 (100%)	1 (0%)	0	100	100
1	C	177/250 (71%)	176 (99%)	1 (1%)	0	100	100
1	D	198/250 (79%)	196 (99%)	2 (1%)	0	100	100
All	All	760/1000 (76%)	755 (99%)	5 (1%)	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	148/207 (72%)	146 (99%)	2 (1%)	67	47
1	B	161/207 (78%)	158 (98%)	3 (2%)	57	34
1	C	145/207 (70%)	143 (99%)	2 (1%)	67	47
1	D	154/207 (74%)	152 (99%)	2 (1%)	69	50
All	All	608/828 (73%)	599 (98%)	9 (2%)	65	44

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	THR
1	A	186	ARG
1	B	18	GLN
1	B	111	TYR
1	B	186	ARG
1	C	51	SER
1	C	156	LEU
1	D	25	ILE
1	D	110	ARG

Sometimes 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 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 [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	190/250 (76%)	0.35	13 (6%) 17 16	14, 27, 42, 52	0
1	B	207/250 (82%)	0.49	20 (9%) 7 6	12, 22, 42, 57	0
1	C	187/250 (74%)	0.47	21 (11%) 5 4	13, 23, 43, 50	0
1	D	203/250 (81%)	0.79	31 (15%) 2 1	12, 25, 50, 53	0
All	All	787/1000 (78%)	0.53	85 (10%) 5 5	12, 25, 47, 57	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	25	ILE	7.4
1	D	111	TYR	6.9
1	B	226	LEU	6.4
1	D	56	GLY	6.2
1	B	111	TYR	6.2
1	D	44	VAL	6.1
1	B	178	LEU	5.2
1	B	202	GLY	5.1
1	B	216	LEU	4.9
1	A	0	SER	4.7
1	C	201	SER	4.5
1	C	210	TRP	4.5
1	B	58	GLY	4.4
1	B	42	TYR	4.3
1	D	202	GLY	4.3
1	A	216	LEU	4.2
1	D	42	TYR	4.1
1	A	226	LEU	3.9
1	C	200	LEU	3.9
1	D	28	GLY	3.8
1	C	226	LEU	3.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	227	ALA	3.8
1	D	21	PRO	3.7
1	B	157	THR	3.6
1	D	157	THR	3.6
1	A	30	VAL	3.6
1	A	57	PRO	3.5
1	A	224	PRO	3.5
1	D	24	ALA	3.5
1	D	18	GLN	3.4
1	C	225	ASP	3.3
1	C	44	VAL	3.2
1	B	11	ALA	3.2
1	D	22	GLY	3.2
1	A	200	LEU	3.2
1	D	54	GLY	3.2
1	D	58	GLY	3.2
1	C	15	PRO	3.1
1	B	224	PRO	3.1
1	D	26	ALA	3.1
1	C	202	GLY	3.1
1	D	27	SER	3.0
1	C	111	TYR	3.0
1	A	44	VAL	2.9
1	A	111	TYR	2.9
1	B	225	ASP	2.9
1	C	58	GLY	2.8
1	B	18	GLN	2.8
1	B	55	GLY	2.8
1	B	56	GLY	2.8
1	D	55	GLY	2.8
1	D	145	ALA	2.7
1	D	178	LEU	2.7
1	B	14	ASP	2.6
1	C	46	HIS	2.6
1	B	147	VAL	2.5
1	C	204	HIS	2.5
1	D	146	VAL	2.5
1	D	106	PHE	2.5
1	D	149	ILE	2.5
1	A	16	LEU	2.4
1	A	156	LEU	2.4
1	D	223	ARG	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	45	HIS	2.4
1	D	219	THR	2.4
1	D	179	LEU	2.4
1	C	144	ALA	2.4
1	C	203	ASP	2.3
1	C	106	PHE	2.3
1	A	46	HIS	2.3
1	C	181	GLY	2.3
1	C	198	ILE	2.2
1	D	29	LEU	2.2
1	C	205	ALA	2.2
1	A	181	GLY	2.2
1	D	221	GLU	2.1
1	C	107	ALA	2.1
1	B	106	PHE	2.1
1	D	5	VAL	2.1
1	D	6	VAL	2.1
1	D	220	ARG	2.1
1	B	205	ALA	2.1
1	B	17	ARG	2.0
1	B	81	VAL	2.0
1	D	11	ALA	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.