

wwPDB X-ray Structure Validation Summary Report (i)

Nov 14, 2023 – 10:26 PM JST

PDB ID	:	6AIE
Title	:	Crystal structure of a new form of RsmD-like RNA methyl transferase from
		Mycobacterium tuberculosis determined at 1.74 A resolution
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Deposited on		
Resolution	:	1.74 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 (i) 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 (1)) were used in the production of this report:

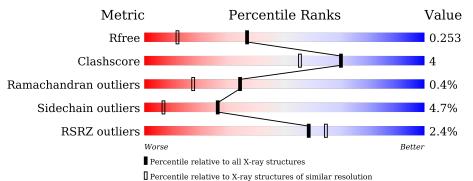
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.74 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	$\begin{array}{l} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	3764(1.76-1.72)
Clashscore	141614	3923 (1.76-1.72)
Ramachandran outliers	138981	3878 (1.76-1.72)
Sidechain outliers	138945	3878 (1.76-1.72)
RSRZ outliers	127900	3705 (1.76-1.72)

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 <=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	А	189	83%	12%	•	·
1	В	189	2% 84%	11%	•	·
1	С	189	3%	9%	•	•



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4354 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.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Δ	182	Total	С	Ν	Ο	S	0	0	0
1		162	1337	830	252	253	2	0	0	0
1	В	182	Total	С	Ν	Ο	S	0	0	0
	D	162	1337	830	252	253	2	0	0	0
1	С	182	Total	С	Ν	Ο	S	0	0	0
		102	1337	830	252	253	2		U	0

• Molecule 1 is a protein called Putative methyltransferase.

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	0	SER	-	expression tag	UNP A0A045IDU3
В	0	SER	-	expression tag	UNP A0A045IDU3
С	0	SER	-	expression tag	UNP A0A045IDU3

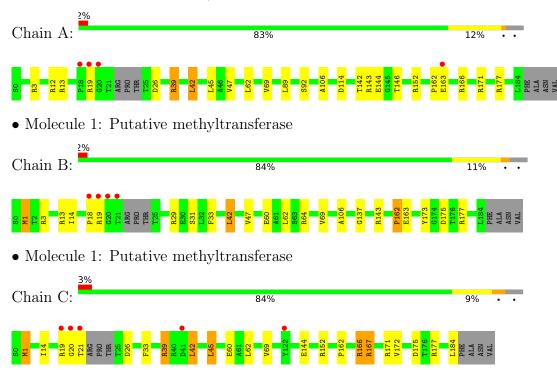
• Molecule 2 is water.

Mol	Chain	Residues	tesidues Atoms		AltConf
2	А	111	Total O 111 111	0	0
2	В	119	Total O 119 119	0	0
2	С	113	Total O 113 113	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: Putative methyltransferase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	95.78Å 55.30 Å 100.74 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.39 - 1.74	Depositor
Resolution (A)	50.37 - 1.74	EDS
% Data completeness	97.8 (50.39-1.74)	Depositor
(in resolution range)	97.8(50.37-1.74)	EDS
R _{merge}	(Not available)	Depositor
$\frac{\mathbf{R}_{sym}}{< I/\sigma(I) > 1}$	0.15	Depositor
$< I/\sigma(I) > 1$	1.73 (at 1.74 Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
D D	0.211 , 0.243	Depositor
R, R_{free}	0.222 , 0.253	DCC
R_{free} test set	1213 reflections (2.27%)	wwPDB-VP
Wilson B-factor $(Å^2)$	17.3	Xtriage
Anisotropy	0.651	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 31.0	EDS
L-test for $twinning^2$	$< L > = 0.49, < L^2 > = 0.32$	Xtriage
	0.034 for -1/2*h-3/2*k,-1/2*h+1/2*k,-1	
	0.033 for -1/2*h+3/2*k,1/2*h+1/2*k,-l	
Estimated twinning fraction	0.438 for $1/2$ *h- $3/2$ *k,- $1/2$ *h- $1/2$ *k,-l	Xtriage
	0.459 for 1/2 *h + 3/2 *k, 1/2 *h - 1/2 *k, -1	
	0.032 for -h,-k,l	
F_o, F_c correlation	0.94	EDS
Total number of atoms	4354	wwPDB-VP
Average B, all atoms $(Å^2)$	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.59% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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		
	Ullalli	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.90	0/1355	1.12	10/1847~(0.5%)	
1	В	0.92	0/1355	0.99	3/1847~(0.2%)	
1	С	0.93	0/1355	1.03	4/1847~(0.2%)	
All	All	0.92	0/4065	1.05	17/5541~(0.3%)	

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	1
1	В	0	2
1	С	0	1
All	All	0	4

There are no bond length outliers.

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	171	ARG	NE-CZ-NH1	9.13	124.87	120.30
1	А	177	ARG	NE-CZ-NH1	9.07	124.83	120.30
1	С	177	ARG	NE-CZ-NH1	8.18	124.39	120.30
1	А	13	ARG	NE-CZ-NH1	8.03	124.32	120.30
1	А	39	ARG	NE-CZ-NH1	-7.88	116.36	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	162	PRO	Peptide
1	В	162	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	В	18	PRO	Peptide
1	С	162	PRO	Peptide

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	А	1337	0	1362	6	0
1	В	1337	0	1362	14	0
1	С	1337	0	1362	10	0
2	А	111	0	0	0	0
2	В	119	0	0	2	0
2	С	113	0	0	3	0
All	All	4354	0	4086	29	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 4.

The worst 5 of 29 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:175:ASP:HA	2:B:220:HOH:O	1.86	0.76
1:B:3:ARG:O	1:B:29:ARG:NH2	2.24	0.71
1:C:167:ARG:NH1	2:C:201:HOH:O	2.26	0.67
1:C:62:LEU:HD13	1:C:69:VAL:HG23	1.83	0.61
1:C:175:ASP:HB3	2:C:310:HOH:O	2.01	0.59

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.



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
1	А	178/189~(94%)	170 (96%)	7~(4%)	1 (1%)	25	10
1	В	178/189~(94%)	172 (97%)	6 (3%)	0	100	100
1	С	178/189~(94%)	172 (97%)	5(3%)	1 (1%)	25	10
All	All	534/567~(94%)	514 (96%)	18 (3%)	2(0%)	34	17

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	19	ARG
1	С	20	GLY

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	134/141~(95%)	126~(94%)	8 (6%)	19 4
1	В	134/141~(95%)	131 (98%)	3~(2%)	52 29
1	С	134/141~(95%)	126 (94%)	8 (6%)	19 4
All	All	402/423~(95%)	383~(95%)	19 (5%)	26 7

5 of 19 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	26	ASP
1	С	45	LEU
1	С	171	ARG
1	С	42	LEU
1	В	1	MET

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. There are no such side chains 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, 95^{th} 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 $>$	#RSR2	$\mathbf{z}{>}2$	$OWAB(Å^2)$	Q < 0.9
1	А	182/189~(96%)	-0.08	4 (2%) 62	68	13, 21, 41, 100	0
1	В	182/189~(96%)	-0.08	4 (2%) 62	68	12, 21, 42, 100	0
1	С	182/189~(96%)	-0.08	5 (2%) 54	60	11, 21, 42, 108	0
All	All	546/567~(96%)	-0.08	13 (2%) 59) 65	11, 21, 42, 108	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	20	GLY	7.9
1	С	19	ARG	6.7
1	В	19	ARG	6.3
1	А	20	GLY	5.0
1	В	21	THR	4.4

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.

