



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 13, 2017 – 10:19 am GMT

PDB ID : 4JIS
Title : Crystal structure of ribitol 5-phosphate cytidylyltransferase (TarI) from *Bacillus subtilis*
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Deposited on : 2013-03-06
Resolution : 1.77 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

<http://wwpdb.org/validation/2016/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) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

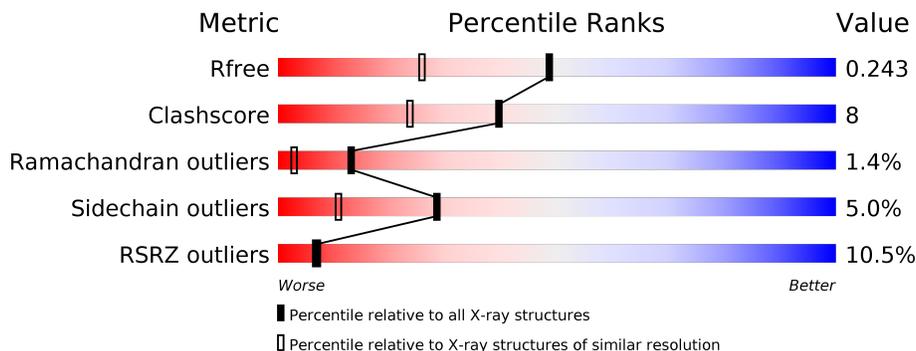
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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}	100719	7172 (1.80-1.76)
Clashscore	112137	8247 (1.80-1.76)
Ramachandran outliers	110173	8154 (1.80-1.76)
Sidechain outliers	110143	8153 (1.80-1.76)
RSRZ outliers	101464	7262 (1.80-1.76)

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. 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	247	
1	B	247	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 3841 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 ribitol-5-phosphate cytidyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	221	1766	1130	296	334	6	0	0	0
1	B	219	1750	1119	294	332	5	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	EXPRESSION TAG	UNP Q8RKI9
A	0	GLY	-	EXPRESSION TAG	UNP Q8RKI9
A	238	LEU	-	EXPRESSION TAG	UNP Q8RKI9
A	239	GLU	-	EXPRESSION TAG	UNP Q8RKI9
A	240	HIS	-	EXPRESSION TAG	UNP Q8RKI9
A	241	HIS	-	EXPRESSION TAG	UNP Q8RKI9
A	242	HIS	-	EXPRESSION TAG	UNP Q8RKI9
A	243	HIS	-	EXPRESSION TAG	UNP Q8RKI9
A	244	HIS	-	EXPRESSION TAG	UNP Q8RKI9
A	245	HIS	-	EXPRESSION TAG	UNP Q8RKI9
B	-1	MET	-	EXPRESSION TAG	UNP Q8RKI9
B	0	GLY	-	EXPRESSION TAG	UNP Q8RKI9
B	238	LEU	-	EXPRESSION TAG	UNP Q8RKI9
B	239	GLU	-	EXPRESSION TAG	UNP Q8RKI9
B	240	HIS	-	EXPRESSION TAG	UNP Q8RKI9
B	241	HIS	-	EXPRESSION TAG	UNP Q8RKI9
B	242	HIS	-	EXPRESSION TAG	UNP Q8RKI9
B	243	HIS	-	EXPRESSION TAG	UNP Q8RKI9
B	244	HIS	-	EXPRESSION TAG	UNP Q8RKI9
B	245	HIS	-	EXPRESSION TAG	UNP Q8RKI9

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	160	Total 160	O 160	0	0
2	B	165	Total 165	O 165	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	103.75Å 60.98Å 91.81Å 90.00° 113.44° 90.00°	Depositor
Resolution (Å)	24.71 – 1.77 24.71 – 1.77	Depositor EDS
% Data completeness (in resolution range)	97.4 (24.71-1.77) 97.4 (24.71-1.77)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.19 (at 1.77Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.206 , 0.241 0.208 , 0.243	Depositor DCC
R_{free} test set	2533 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	25.6	Xtrriage
Anisotropy	0.018	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.015 for $1/2^*h+3/2^*k, 1/2^*h-1/2^*k, -1/2^*h-1/2^*k-l$ 0.011 for $1/2^*h-3/2^*k, -1/2^*h-1/2^*k, -1/2^*h+1/2^*k-l$	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3841	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.23% 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	0.39	0/1793	0.61	2/2423 (0.1%)
1	B	0.39	0/1777	0.64	0/2401
All	All	0.39	0/3570	0.62	2/4824 (0.0%)

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	A	0	2
1	B	0	1
All	All	0	3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	168	THR	C-N-CD	-6.68	105.90	120.60
1	A	117	LEU	CA-CB-CG	5.13	127.11	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	168	THR	Mainchain,Peptide
1	B	160	ARG	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	A	1766	0	1813	34	0
1	B	1750	0	1794	31	0
2	A	160	0	0	5	4
2	B	165	0	0	8	4
All	All	3841	0	3607	60	4

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 8.

The worst 5 of 60 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:205:LYS:NZ	2:B:448:HOH:O	1.97	0.94
1:A:111:ASP:H	1:A:168:THR:HG21	1.39	0.88
1:A:159:VAL:O	2:A:376:HOH:O	1.96	0.84
1:B:111:ASP:H	1:B:168:THR:HG21	1.44	0.82
1:B:152:GLU:OE1	2:B:404:HOH:O	2.04	0.75

All (4) 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)
2:A:400:HOH:O	2:B:393:HOH:O[3_455]	1.56	0.64
2:A:428:HOH:O	2:B:435:HOH:O[3_455]	1.81	0.39
2:A:397:HOH:O	2:B:331:HOH:O[3_455]	2.01	0.19
2:A:431:HOH:O	2:B:437:HOH:O[3_455]	2.08	0.12

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	217/247 (88%)	205 (94%)	6 (3%)	6 (3%)	6	1
1	B	215/247 (87%)	208 (97%)	7 (3%)	0	100	100
All	All	432/494 (87%)	413 (96%)	13 (3%)	6 (1%)	13	3

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	161	ASP
1	A	162	ASN
1	A	189	LYS
1	A	191	VAL
1	A	187	GLU

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	200/222 (90%)	191 (96%)	9 (4%)	32	13
1	B	198/222 (89%)	187 (94%)	11 (6%)	25	8
All	All	398/444 (90%)	378 (95%)	20 (5%)	28	11

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	7	LEU
1	B	23	GLN
1	B	168	THR
1	A	197	LYS
1	A	225	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 [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	221/247 (89%)	0.75	22 (9%) 8 7	22, 38, 61, 75	0
1	B	219/247 (88%)	0.59	24 (10%) 6 6	21, 37, 56, 70	0
All	All	440/494 (89%)	0.67	46 (10%) 7 7	21, 37, 57, 75	0

The worst 5 of 46 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	226	VAL	8.7
1	A	190	ALA	8.5
1	A	191	VAL	6.7
1	A	161	ASP	6.4
1	A	162	ASN	4.8

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.