

Full wwPDB X-ray Structure Validation Report (i)

May 13, 2020 - 05:03 am BST

PDB ID	:	5DTP
Title	:	Crystal structure of M. tuberculosis EchA6 (apo, trigonal crystal form)
Authors	:	Cox, J.A.G.; Besra, G.S.; Futterer, K.
Deposited on		
Resolution	:	1.91 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

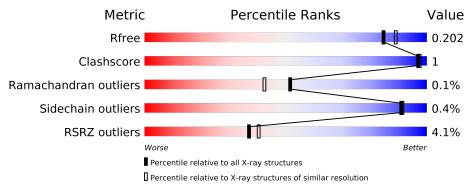
MolProbity		
$\mathbf{Xtriage} \ (\mathbf{Phenix})$:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
$\operatorname{CCP4}$:	$7.0.044 (\mathrm{Gargrove})$
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.91 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
R_{free}	130704	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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 <=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	263	86%	·	10%
1	В	263	4%	•	12%
1	С	263	90%		• 7%



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5610 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		Ate	oms			ZeroOcc	AltConf	Trace
1	Λ	236	Total	С	Ν	Ο	\mathbf{S}	0	0	0
	л	230	1748	1098	316	326	8	0	0	0
1	1 B	232	Total	С	Ν	Ο	S	0	0	0
		202	1725	1082	313	322	8	0	0	0
1	C	244	Total	С	Ν	Ο	S	0	0	0
	$\angle 44$	1811	1137	328	337	9	0	0	U	

• Molecule 1 is a protein called Probable enoyl-CoA hydratase echA6.

Chain	Residue	Modelled	Actual	Comment	Reference
А	-19	MET	-	initiating methionine	UNP P9WNP1
A	-18	GLY	-	expression tag	UNP P9WNP1
A	-17	SER	-	expression tag	UNP P9WNP1
А	-16	SER	-	expression tag	UNP P9WNP1
A	-15	HIS	-	expression tag	UNP P9WNP1
А	-14	HIS	-	expression tag	UNP P9WNP1
A	-13	HIS	-	expression tag	UNP P9WNP1
A	-12	HIS	-	expression tag	UNP P9WNP1
A	-11	HIS	-	expression tag	UNP P9WNP1
A	-10	HIS	-	expression tag	UNP P9WNP1
А	-9	SER	-	expression tag	UNP P9WNP1
A	-8	SER	-	expression tag	UNP P9WNP1
А	-7	GLY	-	expression tag	UNP P9WNP1
A	-6	LEU	-	expression tag	UNP P9WNP1
A	-5	VAL	-	expression tag	UNP P9WNP1
А	-4	PRO	-	expression tag	UNP P9WNP1
A	-3	ARG	-	expression tag	UNP P9WNP1
А	-2	GLY	-	expression tag	UNP P9WNP1
A	-1	SER	-	expression tag	UNP P9WNP1
А	0	ALA	-	expression tag	UNP P9WNP1
В	-19	MET	-	initiating methionine	UNP P9WNP1
В	-18	GLY	-	expression tag	UNP P9WNP1
В	-17	SER	-	expression tag	UNP P9WNP1

There are 60 discrepancies between the modelled and reference sequences:

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Chain	Residue	Modelled	Actual	Comment	Reference
В	-16	SER	-	expression tag	UNP P9WNP1
В	-15	HIS	_	expression tag	UNP P9WNP1
В	-14	HIS	_	expression tag	UNP P9WNP1
В	-13	HIS	_	expression tag	UNP P9WNP1
В	-12	HIS	_	expression tag	UNP P9WNP1
В	-11	HIS	_	expression tag	UNP P9WNP1
В	-10	HIS	-	expression tag	UNP P9WNP1
В	-9	SER	-	expression tag	UNP P9WNP1
В	-8	SER	-	expression tag	UNP P9WNP1
В	-7	GLY	-	expression tag	UNP P9WNP1
В	-6	LEU	-	expression tag	UNP P9WNP1
В	-5	VAL	-	expression tag	UNP P9WNP1
В	-4	PRO	-	expression tag	UNP P9WNP1
В	-3	ARG	-	expression tag	UNP P9WNP1
В	-2	GLY	-	expression tag	UNP P9WNP1
В	-1	SER	-	expression tag	UNP P9WNP1
В	0	ALA	-	expression tag	UNP P9WNP1
С	-19	MET	-	initiating methionine	UNP P9WNP1
С	-18	GLY	-	expression tag	UNP P9WNP1
С	-17	SER	-	expression tag	UNP P9WNP1
С	-16	SER	-	expression tag	UNP P9WNP1
С	-15	HIS	_	expression tag	UNP P9WNP1
С	-14	HIS	-	expression tag	UNP P9WNP1
С	-13	HIS	_	expression tag	UNP P9WNP1
С	-12	HIS	-	expression tag	UNP P9WNP1
C	-11	HIS	_	expression tag	UNP P9WNP1
С	-10	HIS	-	expression tag	UNP P9WNP1
С	-9	SER	-	expression tag	UNP P9WNP1
С	-8	SER	-	expression tag	UNP P9WNP1
C	-7	GLY	-	expression tag	UNP P9WNP1
C	-6	LEU	-	expression tag	UNP P9WNP1
С	-5	VAL	-	expression tag	UNP P9WNP1
C	-4	PRO	-	expression tag	UNP P9WNP1
С	-3	ARG	-	expression tag	UNP P9WNP1
C	-2	GLY	-	expression tag	UNP P9WNP1
С	-1	SER	-	expression tag	UNP P9WNP1
C	0	ALA	-	expression tag	UNP P9WNP1

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• Molecule 2 is water.

$ \begin{array}{ c c c c c c c c c c c c c c c c c c c$	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	2	А	96		0	0

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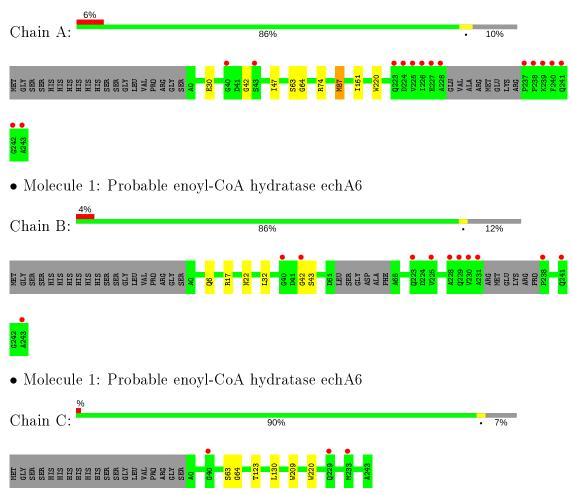
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	115	Total O 115 115	0	0
2	С	115	Total O 115 115	0	0



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: Probable enoyl-CoA hydratase echA6





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants	103.12Å 103.12 Å 143.44 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	89.30 - 1.91	Depositor
Resolution (A)	89.30 - 1.91	EDS
% Data completeness	$100.0 \ (89.30-1.91)$	Depositor
(in resolution range)	$100.0 \ (89.30-1.91)$	EDS
R _{merge}	0.07	Depositor
R _{sym}	0.07	Depositor
$< I/\sigma(I) > 1$	$2.72 (at 1.91 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
D D .	0.182 , 0.201	Depositor
R, R_{free}	0.181 , 0.202	DCC
R_{free} test set	3508 reflections $(5.09%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	28.9	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.35 , 41.1	EDS
L-test for twinning ²	$< L > = 0.48, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5610	wwPDB-VP
Average B, all atoms $(Å^2)$	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.00% 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	Boı	nd lengths	Bond angles		
	Cham	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.47	1/1777~(0.1%)	0.51	0/2409	
1	В	0.47	0/1753	0.50	0/2376	
1	С	0.48	2/1843~(0.1%)	0.51	0/2501	
All	All	0.47	3/5373~(0.1%)	0.51	0/7286	

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
1	С	209	TRP	CD2-CE2	5.11	1.47	1.41
1	С	220	TRP	CD2-CE2	5.10	1.47	1.41
1	А	220	TRP	CD2-CE2	5.04	1.47	1.41

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	А	1748	0	1735	4	0
1	В	1725	0	1722	3	0
1	С	1811	0	1801	2	0
2	А	96	0	0	1	0
2	В	115	0	0	0	0
2	С	115	0	0	0	0
All	All	5610	0	5258	9	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 (9) 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:161:ILE:HD12	2:A:369:HOH:O	2.15	0.46
1:A:63:SER:HA	1:A:64:GLY:HA2	1.75	0.45
1:B:42:GLY:HA2	1:B:43:SER:OG	2.17	0.45
1:A:30:GLU:HB2	1:A:74:ARG:HD3	2.00	0.44
1:C:63:SER:HA	1:C:64:GLY:HA2	1.77	0.43
1:B:17:ARG:HB2	1:B:22:ASN:HA	2.01	0.41
1:C:123:THR:HG21	1:C:130:LEU:HG	2.00	0.41
1:A:47:ILE:HG13	1:A:87:MET:HE2	2.02	0.41
1:B:6:GLN:OE1	1:B:43:SER:HB2	2.21	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	А	232/263~(88%)	222~(96%)	9~(4%)	1 (0%)	34 24
1	В	226/263~(86%)	216~(96%)	10 (4%)	0	100 100
1	С	242/263~(92%)	236~(98%)	6(2%)	0	100 100
All	All	700/789~(89%)	674 (96%)	25~(4%)	1 (0%)	51 42

All (1) Ramachandran outliers are listed below:

Mol	Chain	\mathbf{Res}	Type
1	А	42	GLY



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	А	167/196~(85%)	166~(99%)	1 (1%)	86 86
1	В	167/196~(85%)	166~(99%)	1 (1%)	86 86
1	С	174/196~(89%)	174 (100%)	0	100 100
All	All	508/588~(86%)	506 (100%)	2(0%)	91 91

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	87	MET
1	В	32	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	6	GLN
1	А	178	GLN
1	В	178	GLN
1	В	223	GLN
1	С	6	GLN
1	С	178	GLN

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, 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 $>$	# RSRZ > 2	$OWAB(Å^2)$	$Q{<}0.9$
1	А	236/263~(89%)	0.03	15 (6%) 19 22	19, 31, 63, 107	0
1	В	232/263~(88%)	-0.04	11 (4%) 31 34	18, 29, 60, 85	0
1	С	244/263~(92%)	-0.18	3 (1%) 79 81	18, 29, 52, 66	0
All	All	712/789~(90%)	-0.07	29 (4%) 37 40	18, 30, 57, 107	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	237	PRO	17.1
1	А	238	PRO	7.0
1	А	223	GLN	5.2
1	А	243	ALA	4.6
1	С	233	MET	4.5
1	В	243	ALA	4.4
1	А	225	VAL	4.1
1	А	239	LYS	4.0
1	А	241	GLN	4.0
1	А	40	GLY	3.8
1	В	230	VAL	3.8
1	В	241	GLN	3.5
1	А	224	ASP	3.3
1	В	229	GLN	3.3
1	А	227	GLU	3.2
1	В	231	ALA	3.1
1	В	40	GLY	3.0
1	А	43	SER	3.0
1	В	42	GLY	2.9
1	А	240	PHE	2.7
1	С	229	GLN	2.7
1	В	223	GLN	2.7
1	В	228	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	В	238	PRO	2.6
1	А	242	GLY	2.6
1	В	225	VAL	2.2
1	С	40	GLY	2.1
1	А	228	ALA	2.1
1	А	226	ILE	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.

