

wwPDB X-ray Structure Validation Summary Report (i)

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PDB ID : 6MEL

Title : Succinyl-CoA synthase from Campylobacter jejuni

Authors: Osipiuk, J.; Maltseva, N.; Jedrzejczak, R.; Satchell, K.J.F.; Joachimiak, A.;

Center for Structural Genomics of Infectious Diseases (CSGID)

Deposited on : 2018-09-06

Resolution : 2.06 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 (i)) were used in the production of this report:

MolProbity : FAILED

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : FAILED

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

Validation Pipeline (wwPDB-VP) : 2.35.1

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 2.06 Å.

There are no overall percentile quality scores available for this entry.

MolProbity and EDS failed to run properly - the sequence quality summary graphics cannot be shown.



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5585 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 Succinate—CoA ligase subunit alpha.

Mol	Chain	Residues		Atoms				ZeroOcc	AltConf	Trace
1	٨	289	Total	С	N	О	S	0	9	0
1	A	209	2117	1356	348	402	11	0		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	290	GLU	-	expression tag	UNP A0A2U0Q9D2
A	291	ASN	-	1	UNP A0A2U0Q9D2
A	292	LEU	-	expression tag	UNP A0A2U0Q9D2
A	293	TYR	-	expression tag	UNP A0A2U0Q9D2
A	294	PHE	_	expression tag	UNP A0A2U0Q9D2
A	295	GLN	-	expression tag	UNP A0A2U0Q9D2

• Molecule 2 is a protein called Succinate—CoA ligase [ADP-forming] subunit beta.

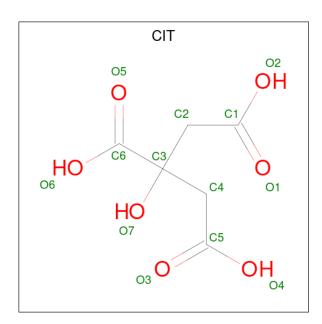
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	В	386	Total 2995	C 1909	N 501	O 570	S 15	0	12	0

• Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0

• Molecule 4 is CITRIC ACID (three-letter code: CIT) (formula: C₆H₈O₇).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	В	1	Total 13	C 6	O 7	0	0

• Molecule 5 is water.

	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
Ī	5	A	200	Total O 200 200	0	0
	5	В	259	Total O 259 259	0	0

MolProbity and EDS failed to run properly - this section is therefore empty.



3 Data and refinement statistics (i)

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants	123.07Å 123.07Å 146.79Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.90 - 2.06	Depositor
% Data completeness	99.9 (48.90-2.06)	Depositor
(in resolution range)	,	
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.72 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.150 , 0.175	Depositor
Wilson B-factor $(Å^2)$	42.3	Xtriage
Anisotropy	0.167	Xtriage
L-test for twinning ²	$< L > = 0.49, < L^2> = 0.32$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
Total number of atoms	5585	wwPDB-VP
Average B, all atoms (\mathring{A}^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.17% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ 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.

4 Model quality (i)

4.1 Standard geometry (i)

MolProbity failed to run properly - this section is therefore empty.

4.2 Too-close contacts (i)

MolProbity failed to run properly - this section is therefore empty.

4.3 Torsion angles (i)

4.3.1 Protein backbone (i)

MolProbity failed to run properly - this section is therefore empty.

4.3.2 Protein sidechains (i)

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA (i)

MolProbity failed to run properly - this section is therefore empty.

4.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

4.5 Carbohydrates (i)

There are no monosaccharides in this entry.

4.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mo	Type	Chain	Pog	Link	Bo	ond leng	Bond angles			
1010	Type	Chain	rtes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	CIT	В	501	-	12,12,12	1.22	0	17,17,17	1.58	3 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

\mathbf{Mol}	Type	Chain	Res	Link	Chirals	Torsions	Rings	
4	CIT	В	501	-	-	2/16/16/16	-	

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
4	В	501	CIT	O6-C6-C3	3.55	119.22	113.05
4	В	501	CIT	O4-C5-C4	2.85	123.51	114.35
4	В	501	CIT	O5-C6-C3	-2.73	118.39	122.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	В	501	CIT	C3-C4-C5-O4
4	В	501	CIT	C3-C4-C5-O3

There are no ring outliers.

No monomer is involved in short contacts.

4.7 Other polymers (i)

There are no such residues in this entry.

4.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



5 Fit of model and data (i)

5.1 Protein, DNA and RNA chains (i)

EDS failed to run properly - this section is therefore empty.

5.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS failed to run properly - this section is therefore empty.

5.3 Carbohydrates (i)

EDS failed to run properly - this section is therefore empty.

5.4 Ligands (i)

EDS failed to run properly - this section is therefore empty.

5.5 Other polymers (i)

EDS failed to run properly - this section is therefore empty.

