

# Full wwPDB X-ray Structure Validation Report (i)

#### Oct 3, 2023 – 06:29 AM EDT

PDB ID	:	6ULM
Title	:	Crystal structure of human cadherin 17 EC1-2
Authors	:	Gray, M.E.; Sotomayor, M.
Deposited on		
Resolution	:	2.15  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 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	:	FAILED
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	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:  $X\hbox{-}RAY\,DIFFRACTION$ 

The reported resolution of this entry is 2.15 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 3 unique types of molecules in this entry. The entry contains 3401 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	Atoms					ZeroOcc	AltConf	Trace
1	В	209	Total	С	Ν	Ο	$\mathbf{S}$	0	2	0
	I D	209	1654	1046	275	328	5			
1	Λ	194	Total	С	Ν	0	S	0	0	0
	A	194	1538	969	256	309	4	0		

• Molecule 1 is a protein called Cadherin-17.

Chain	Residue	Modelled	Actual	Comment	Reference
В	0	MET	-	initiating methionine	UNP Q12864
В	93	GLU	LYS	variant	UNP Q12864
В	220	LEU	-	expression tag	UNP Q12864
В	221	GLU	-	expression tag	UNP Q12864
В	222	HIS	-	expression tag	UNP Q12864
В	223	HIS	-	expression tag	UNP Q12864
В	224	HIS	-	expression tag	UNP Q12864
В	225	HIS	-	expression tag	UNP Q12864
В	226	HIS	-	expression tag	UNP Q12864
В	227	HIS	-	expression tag	UNP Q12864
А	0	MET	-	initiating methionine	UNP Q12864
А	93	GLU	LYS	variant	UNP Q12864
А	220	LEU	-	expression tag	UNP Q12864
А	221	GLU	-	expression tag	UNP Q12864
A	222	HIS	-	expression tag	UNP Q12864
А	223	HIS	-	expression tag	UNP Q12864
А	224	HIS	-	expression tag	UNP Q12864
А	225	HIS	-	expression tag	UNP Q12864
А	226	HIS	-	expression tag	UNP Q12864
А	227	HIS	-	expression tag	UNP Q12864

There are 20 discrepancies between the modelled and reference sequences:

• Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	3	Total Ca 3 3	0	0
2	А	3	Total Ca 3 3	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	В	115	Total O 115 115	0	0
3	А	88	Total O   88 88	0	0

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



# 3 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	58.39Å 81.61Å 106.81Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	47.53 - 2.15	Depositor
% Data completeness	99.5 (47.53-2.15)	Depositor
(in resolution range)		Depositor
R <sub>merge</sub>	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.54$ (at $2.16\text{\AA}$ )	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
$R, R_{free}$	0.206 , $0.257$	Depositor
Wilson B-factor $(Å^2)$	23.4	Xtriage
Anisotropy	0.801	Xtriage
L-test for twinning <sup>2</sup>	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3401	wwPDB-VP
Average B, all atoms $(Å^2)$	32.0	wwPDB-VP

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

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.88% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<sup>&</sup>lt;sup>1</sup>Intensities 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 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.



There are no torsion outliers.

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.

