



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 2, 2023 – 12:16 AM EDT

PDB ID : 6MX2  
Title : Crystal Structure of ClpP1 from Clostridium difficile 630.  
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Deposited on : 2018-10-30  
Resolution : 2.50 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](mailto: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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) 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

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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

There are 4 unique types of molecules in this entry. The entry contains 20432 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 ATP-dependent Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	182	Total	C	N	O	S	0	0	0
			1403	891	236	268	8			
1	B	185	Total	C	N	O	S	0	0	0
			1425	905	239	273	8			
1	C	181	Total	C	N	O	S	0	0	0
			1395	888	232	267	8			
1	D	181	Total	C	N	O	S	0	0	0
			1393	887	232	266	8			
1	E	182	Total	C	N	O	S	0	0	0
			1399	890	233	268	8			
1	F	183	Total	C	N	O	S	0	0	0
			1407	896	234	269	8			
1	G	183	Total	C	N	O	S	0	0	0
			1408	896	234	270	8			
1	H	181	Total	C	N	O	S	0	0	0
			1394	888	232	266	8			
1	I	180	Total	C	N	O	S	0	0	0
			1385	881	231	265	8			
1	J	187	Total	C	N	O	S	0	0	0
			1441	915	242	276	8			
1	K	182	Total	C	N	O	S	0	0	0
			1400	890	233	269	8			
1	L	183	Total	C	N	O	S	0	0	0
			1407	896	234	269	8			
1	M	181	Total	C	N	O	S	0	0	0
			1392	885	232	267	8			
1	N	180	Total	C	N	O	S	0	0	0
			1378	877	227	266	8			

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Na 1	0	0
2	B	1	Total 1	Na 1	0	0
2	C	1	Total 1	Na 1	0	0
2	D	1	Total 1	Na 1	0	0
2	E	1	Total 1	Na 1	0	0
2	F	1	Total 1	Na 1	0	0
2	G	2	Total 2	Na 2	0	0
2	H	1	Total 1	Na 1	0	0
2	I	2	Total 2	Na 2	0	0
2	J	1	Total 1	Na 1	0	0
2	K	1	Total 1	Na 1	0	0
2	L	2	Total 2	Na 2	0	0
2	M	1	Total 1	Na 1	0	0
2	N	1	Total 1	Na 1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	M	1	Total	C	O	0	0
			6	3	3		
3	N	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	55	Total	O	0	0
			55	55		
4	B	47	Total	O	0	0
			47	47		
4	C	62	Total	O	0	0
			62	62		
4	D	60	Total	O	0	0
			60	60		
4	E	70	Total	O	0	0
			70	70		
4	F	53	Total	O	0	0
			53	53		
4	G	52	Total	O	0	0
			52	52		
4	H	57	Total	O	0	0
			57	57		
4	I	57	Total	O	0	0
			57	57		
4	J	50	Total	O	0	0
			50	50		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	K	76	Total 76	O 76	0	0
4	L	47	Total 47	O 47	0	0
4	M	41	Total 41	O 41	0	0
4	N	49	Total 49	O 49	0	0

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

### 3 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.33Å 97.37Å 106.49Å 113.15° 104.56° 103.21°	Depositor
Resolution (Å)	34.90 – 2.50	Depositor
% Data completeness (in resolution range)	98.0 (34.90-2.50)	Depositor
$R_{merge}$	0.22	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.84 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, $R_{free}$	0.193 , 0.244	Depositor
Wilson B-factor (Å <sup>2</sup> )	27.3	Xtriage
Anisotropy	0.011	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	20432	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.23% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<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.

## 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 19 ligands modelled in this entry, 17 are monoatomic - leaving 2 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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	GOL	N	201	-	5,5,5	0.94	0	5,5,5	1.04	0
3	GOL	M	201	-	5,5,5	0.90	0	5,5,5	1.00	0

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.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	N	201	-	-	0/4/4/4	-
3	GOL	M	201	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	M	201	GOL	O1-C1-C2-C3
3	M	201	GOL	C1-C2-C3-O3
3	M	201	GOL	O1-C1-C2-O2
3	M	201	GOL	O2-C2-C3-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.