



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 19, 2023 – 02:23 AM EDT

PDB ID : 5DT5  
Title : Crystal structure of the GH1 beta-glucosidase from *Exiguobacterium antarcticum* B7 in space group P21  
Authors : Zanphorlin, L.M.; Giuseppe, P.O.; Tonoli, C.C.C.; Murakami, M.T.  
Deposited on : 2015-09-17  
Resolution : 2.24 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](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 : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35.1

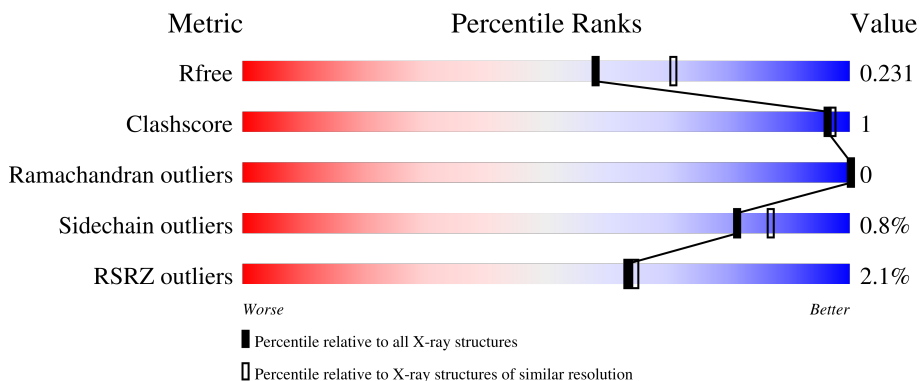
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.24 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.



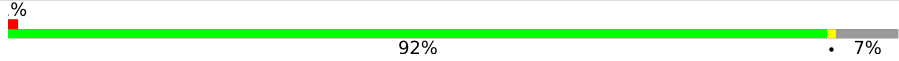

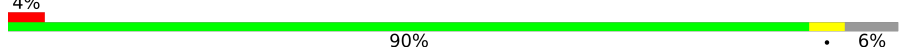
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	2391 (2.26-2.22)
Clashscore	141614	2539 (2.26-2.22)
Ramachandran outliers	138981	2489 (2.26-2.22)
Sidechain outliers	138945	2490 (2.26-2.22)
RSRZ outliers	127900	2353 (2.26-2.22)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ . 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	471	 4% 92% 6%
1	B	471	 % 91% 6%
1	C	471	 % 92% 6%
1	D	471	 % 93% 6%
1	E	471	 2% 90% 7%

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Mol	Chain	Length	Quality of chain
1	F	471	 <p>% 92% 7%</p>
1	G	471	 <p>3% 90% 6%</p>
1	H	471	 <p>4% 90% 6%</p>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 28918 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 Beta-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	444	Total 3613	C 2320	N 603	O 675	S 15	0	0	0
1	B	441	Total 3596	C 2308	N 600	O 672	S 16	0	0	0
1	C	443	Total 3604	C 2314	N 601	O 674	S 15	0	0	0
1	D	443	Total 3604	C 2314	N 601	O 674	S 15	0	0	0
1	E	438	Total 3570	C 2291	N 595	O 669	S 15	0	0	0
1	F	439	Total 3579	C 2297	N 597	O 670	S 15	0	0	0
1	G	441	Total 3596	C 2308	N 600	O 672	S 16	0	0	0
1	H	443	Total 3604	C 2314	N 601	O 674	S 15	0	0	0

There are 184 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	initiating methionine	UNP K0A8J9
A	-21	GLY	-	expression tag	UNP K0A8J9
A	-20	SER	-	expression tag	UNP K0A8J9
A	-19	SER	-	expression tag	UNP K0A8J9
A	-18	HIS	-	expression tag	UNP K0A8J9
A	-17	HIS	-	expression tag	UNP K0A8J9
A	-16	HIS	-	expression tag	UNP K0A8J9
A	-15	HIS	-	expression tag	UNP K0A8J9
A	-14	HIS	-	expression tag	UNP K0A8J9
A	-13	HIS	-	expression tag	UNP K0A8J9
A	-12	SER	-	expression tag	UNP K0A8J9
A	-11	SER	-	expression tag	UNP K0A8J9
A	-10	GLY	-	expression tag	UNP K0A8J9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	LEU	-	expression tag	UNP K0A8J9
A	-8	VAL	-	expression tag	UNP K0A8J9
A	-7	PRO	-	expression tag	UNP K0A8J9
A	-6	ARG	-	expression tag	UNP K0A8J9
A	-5	GLY	-	expression tag	UNP K0A8J9
A	-4	SER	-	expression tag	UNP K0A8J9
A	-3	HIS	-	expression tag	UNP K0A8J9
A	-2	MET	-	expression tag	UNP K0A8J9
A	-1	ALA	-	expression tag	UNP K0A8J9
A	0	SER	-	expression tag	UNP K0A8J9
B	-22	MET	-	initiating methionine	UNP K0A8J9
B	-21	GLY	-	expression tag	UNP K0A8J9
B	-20	SER	-	expression tag	UNP K0A8J9
B	-19	SER	-	expression tag	UNP K0A8J9
B	-18	HIS	-	expression tag	UNP K0A8J9
B	-17	HIS	-	expression tag	UNP K0A8J9
B	-16	HIS	-	expression tag	UNP K0A8J9
B	-15	HIS	-	expression tag	UNP K0A8J9
B	-14	HIS	-	expression tag	UNP K0A8J9
B	-13	HIS	-	expression tag	UNP K0A8J9
B	-12	SER	-	expression tag	UNP K0A8J9
B	-11	SER	-	expression tag	UNP K0A8J9
B	-10	GLY	-	expression tag	UNP K0A8J9
B	-9	LEU	-	expression tag	UNP K0A8J9
B	-8	VAL	-	expression tag	UNP K0A8J9
B	-7	PRO	-	expression tag	UNP K0A8J9
B	-6	ARG	-	expression tag	UNP K0A8J9
B	-5	GLY	-	expression tag	UNP K0A8J9
B	-4	SER	-	expression tag	UNP K0A8J9
B	-3	HIS	-	expression tag	UNP K0A8J9
B	-2	MET	-	expression tag	UNP K0A8J9
B	-1	ALA	-	expression tag	UNP K0A8J9
B	0	SER	-	expression tag	UNP K0A8J9
C	-22	MET	-	initiating methionine	UNP K0A8J9
C	-21	GLY	-	expression tag	UNP K0A8J9
C	-20	SER	-	expression tag	UNP K0A8J9
C	-19	SER	-	expression tag	UNP K0A8J9
C	-18	HIS	-	expression tag	UNP K0A8J9
C	-17	HIS	-	expression tag	UNP K0A8J9
C	-16	HIS	-	expression tag	UNP K0A8J9
C	-15	HIS	-	expression tag	UNP K0A8J9
C	-14	HIS	-	expression tag	UNP K0A8J9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-13	HIS	-	expression tag	UNP K0A8J9
C	-12	SER	-	expression tag	UNP K0A8J9
C	-11	SER	-	expression tag	UNP K0A8J9
C	-10	GLY	-	expression tag	UNP K0A8J9
C	-9	LEU	-	expression tag	UNP K0A8J9
C	-8	VAL	-	expression tag	UNP K0A8J9
C	-7	PRO	-	expression tag	UNP K0A8J9
C	-6	ARG	-	expression tag	UNP K0A8J9
C	-5	GLY	-	expression tag	UNP K0A8J9
C	-4	SER	-	expression tag	UNP K0A8J9
C	-3	HIS	-	expression tag	UNP K0A8J9
C	-2	MET	-	expression tag	UNP K0A8J9
C	-1	ALA	-	expression tag	UNP K0A8J9
C	0	SER	-	expression tag	UNP K0A8J9
D	-22	MET	-	initiating methionine	UNP K0A8J9
D	-21	GLY	-	expression tag	UNP K0A8J9
D	-20	SER	-	expression tag	UNP K0A8J9
D	-19	SER	-	expression tag	UNP K0A8J9
D	-18	HIS	-	expression tag	UNP K0A8J9
D	-17	HIS	-	expression tag	UNP K0A8J9
D	-16	HIS	-	expression tag	UNP K0A8J9
D	-15	HIS	-	expression tag	UNP K0A8J9
D	-14	HIS	-	expression tag	UNP K0A8J9
D	-13	HIS	-	expression tag	UNP K0A8J9
D	-12	SER	-	expression tag	UNP K0A8J9
D	-11	SER	-	expression tag	UNP K0A8J9
D	-10	GLY	-	expression tag	UNP K0A8J9
D	-9	LEU	-	expression tag	UNP K0A8J9
D	-8	VAL	-	expression tag	UNP K0A8J9
D	-7	PRO	-	expression tag	UNP K0A8J9
D	-6	ARG	-	expression tag	UNP K0A8J9
D	-5	GLY	-	expression tag	UNP K0A8J9
D	-4	SER	-	expression tag	UNP K0A8J9
D	-3	HIS	-	expression tag	UNP K0A8J9
D	-2	MET	-	expression tag	UNP K0A8J9
D	-1	ALA	-	expression tag	UNP K0A8J9
D	0	SER	-	expression tag	UNP K0A8J9
E	-22	MET	-	initiating methionine	UNP K0A8J9
E	-21	GLY	-	expression tag	UNP K0A8J9
E	-20	SER	-	expression tag	UNP K0A8J9
E	-19	SER	-	expression tag	UNP K0A8J9
E	-18	HIS	-	expression tag	UNP K0A8J9

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-17	HIS	-	expression tag	UNP K0A8J9
E	-16	HIS	-	expression tag	UNP K0A8J9
E	-15	HIS	-	expression tag	UNP K0A8J9
E	-14	HIS	-	expression tag	UNP K0A8J9
E	-13	HIS	-	expression tag	UNP K0A8J9
E	-12	SER	-	expression tag	UNP K0A8J9
E	-11	SER	-	expression tag	UNP K0A8J9
E	-10	GLY	-	expression tag	UNP K0A8J9
E	-9	LEU	-	expression tag	UNP K0A8J9
E	-8	VAL	-	expression tag	UNP K0A8J9
E	-7	PRO	-	expression tag	UNP K0A8J9
E	-6	ARG	-	expression tag	UNP K0A8J9
E	-5	GLY	-	expression tag	UNP K0A8J9
E	-4	SER	-	expression tag	UNP K0A8J9
E	-3	HIS	-	expression tag	UNP K0A8J9
E	-2	MET	-	expression tag	UNP K0A8J9
E	-1	ALA	-	expression tag	UNP K0A8J9
E	0	SER	-	expression tag	UNP K0A8J9
F	-22	MET	-	initiating methionine	UNP K0A8J9
F	-21	GLY	-	expression tag	UNP K0A8J9
F	-20	SER	-	expression tag	UNP K0A8J9
F	-19	SER	-	expression tag	UNP K0A8J9
F	-18	HIS	-	expression tag	UNP K0A8J9
F	-17	HIS	-	expression tag	UNP K0A8J9
F	-16	HIS	-	expression tag	UNP K0A8J9
F	-15	HIS	-	expression tag	UNP K0A8J9
F	-14	HIS	-	expression tag	UNP K0A8J9
F	-13	HIS	-	expression tag	UNP K0A8J9
F	-12	SER	-	expression tag	UNP K0A8J9
F	-11	SER	-	expression tag	UNP K0A8J9
F	-10	GLY	-	expression tag	UNP K0A8J9
F	-9	LEU	-	expression tag	UNP K0A8J9
F	-8	VAL	-	expression tag	UNP K0A8J9
F	-7	PRO	-	expression tag	UNP K0A8J9
F	-6	ARG	-	expression tag	UNP K0A8J9
F	-5	GLY	-	expression tag	UNP K0A8J9
F	-4	SER	-	expression tag	UNP K0A8J9
F	-3	HIS	-	expression tag	UNP K0A8J9
F	-2	MET	-	expression tag	UNP K0A8J9
F	-1	ALA	-	expression tag	UNP K0A8J9
F	0	SER	-	expression tag	UNP K0A8J9
G	-22	MET	-	initiating methionine	UNP K0A8J9

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-21	GLY	-	expression tag	UNP K0A8J9
G	-20	SER	-	expression tag	UNP K0A8J9
G	-19	SER	-	expression tag	UNP K0A8J9
G	-18	HIS	-	expression tag	UNP K0A8J9
G	-17	HIS	-	expression tag	UNP K0A8J9
G	-16	HIS	-	expression tag	UNP K0A8J9
G	-15	HIS	-	expression tag	UNP K0A8J9
G	-14	HIS	-	expression tag	UNP K0A8J9
G	-13	HIS	-	expression tag	UNP K0A8J9
G	-12	SER	-	expression tag	UNP K0A8J9
G	-11	SER	-	expression tag	UNP K0A8J9
G	-10	GLY	-	expression tag	UNP K0A8J9
G	-9	LEU	-	expression tag	UNP K0A8J9
G	-8	VAL	-	expression tag	UNP K0A8J9
G	-7	PRO	-	expression tag	UNP K0A8J9
G	-6	ARG	-	expression tag	UNP K0A8J9
G	-5	GLY	-	expression tag	UNP K0A8J9
G	-4	SER	-	expression tag	UNP K0A8J9
G	-3	HIS	-	expression tag	UNP K0A8J9
G	-2	MET	-	expression tag	UNP K0A8J9
G	-1	ALA	-	expression tag	UNP K0A8J9
G	0	SER	-	expression tag	UNP K0A8J9
H	-22	MET	-	initiating methionine	UNP K0A8J9
H	-21	GLY	-	expression tag	UNP K0A8J9
H	-20	SER	-	expression tag	UNP K0A8J9
H	-19	SER	-	expression tag	UNP K0A8J9
H	-18	HIS	-	expression tag	UNP K0A8J9
H	-17	HIS	-	expression tag	UNP K0A8J9
H	-16	HIS	-	expression tag	UNP K0A8J9
H	-15	HIS	-	expression tag	UNP K0A8J9
H	-14	HIS	-	expression tag	UNP K0A8J9
H	-13	HIS	-	expression tag	UNP K0A8J9
H	-12	SER	-	expression tag	UNP K0A8J9
H	-11	SER	-	expression tag	UNP K0A8J9
H	-10	GLY	-	expression tag	UNP K0A8J9
H	-9	LEU	-	expression tag	UNP K0A8J9
H	-8	VAL	-	expression tag	UNP K0A8J9
H	-7	PRO	-	expression tag	UNP K0A8J9
H	-6	ARG	-	expression tag	UNP K0A8J9
H	-5	GLY	-	expression tag	UNP K0A8J9
H	-4	SER	-	expression tag	UNP K0A8J9
H	-3	HIS	-	expression tag	UNP K0A8J9

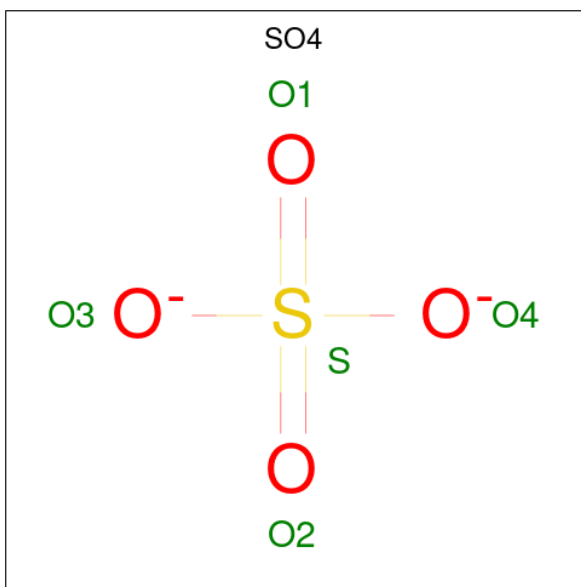
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Chain	Residue	Modelled	Actual	Comment	Reference
H	-2	MET	-	expression tag	UNP K0A8J9
H	-1	ALA	-	expression tag	UNP K0A8J9
H	0	SER	-	expression tag	UNP K0A8J9

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total	O	0	0
			8	8		
3	B	12	Total	O	0	0
			12	12		
3	C	22	Total	O	0	0
			22	22		

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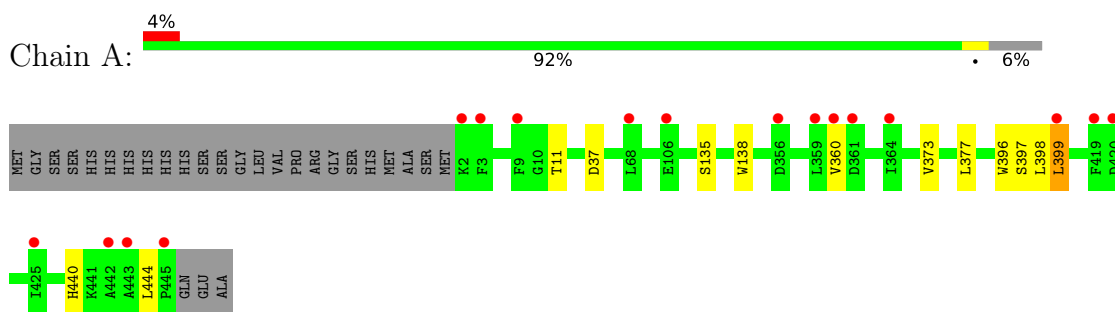
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	50	Total O 50 50	0	0
3	E	11	Total O 11 11	0	0
3	F	14	Total O 14 14	0	0
3	G	4	Total O 4 4	0	0
3	H	6	Total O 6 6	0	0

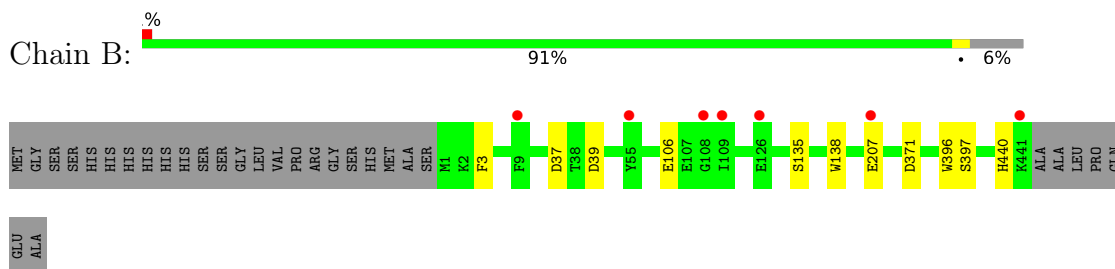
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

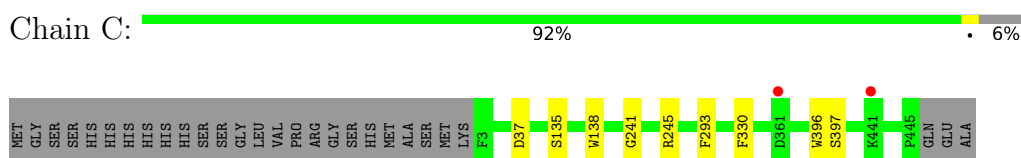
- Molecule 1: Beta-glucosidase



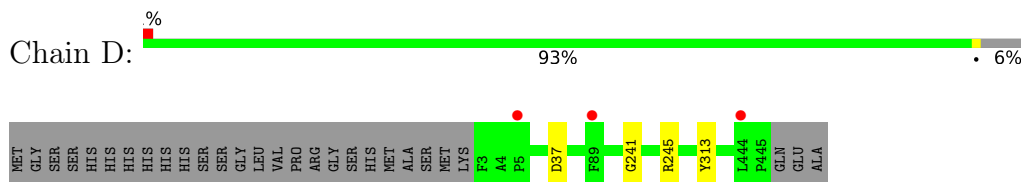
- Molecule 1: Beta-glucosidase



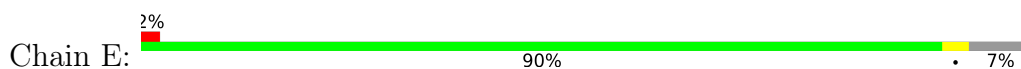
- Molecule 1: Beta-glucosidase



- Molecule 1: Beta-glucosidase



- Molecule 1: Beta-glucosidase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.07Å 104.60Å 199.19Å 90.00° 105.80° 90.00°	Depositor
Resolution (Å)	49.11 – 2.24 49.11 – 2.24	Depositor EDS
% Data completeness (in resolution range)	96.0 (49.11-2.24) 96.0 (49.11-2.24)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.43 (at 2.24Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.209 , 0.229 0.211 , 0.231	Depositor DCC
$R_{free}$ test set	9870 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.2	Xtrriage
Anisotropy	0.387	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 8.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.24$	Xtrriage
Estimated twinning fraction	0.279 for h,-k,-h-l	Xtrriage
Reported twinning fraction	0.722 for H, K, L 0.278 for -H, -K, H+L	Depositor
Outliers	3 of 201050 reflections (0.001%)	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	28918	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.39% 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.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.38	0/3730	0.58	0/5072
1	B	0.39	0/3712	0.59	0/5045
1	C	0.37	0/3721	0.58	0/5061
1	D	0.38	0/3721	0.58	0/5061
1	E	0.38	0/3686	0.58	0/5013
1	F	0.38	0/3695	0.58	0/5024
1	G	0.38	0/3712	0.59	0/5045
1	H	0.39	0/3721	0.60	0/5061
All	All	0.38	0/29698	0.59	0/40382

There are no bond length outliers.

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	A	3613	0	3349	6	0
1	B	3596	0	3333	4	1
1	C	3604	0	3336	4	0
1	D	3604	0	3336	2	0
1	E	3570	0	3295	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	3579	0	3308	3	0
1	G	3596	0	3333	8	0
1	H	3604	0	3336	21	1
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	H	5	0	0	0	0
3	A	8	0	0	0	0
3	B	12	0	0	0	0
3	C	22	0	0	0	0
3	D	50	0	0	0	0
3	E	11	0	0	0	0
3	F	14	0	0	0	0
3	G	4	0	0	0	0
3	H	6	0	0	0	0
All	All	28918	0	26626	53	1

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 (53) 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:H:82:ILE:HD11	1:H:145:VAL:HG11	1.53	0.90
1:H:82:ILE:HG21	1:H:142:PHE:CE1	2.06	0.90
1:E:105:GLN:NE2	1:E:153:ILE:O	2.06	0.89
1:H:89:PHE:HE1	1:H:145:VAL:HG11	1.55	0.72
1:H:89:PHE:CE1	1:H:145:VAL:HG11	2.31	0.64
1:E:101:ALA:HA	1:E:104:LEU:HD12	1.83	0.61
1:E:81:ARG:O	1:E:90:ASN:ND2	2.33	0.60
1:H:144:ARG:HB2	1:H:203:MET:HE1	1.84	0.60
1:H:98:LYS:HE2	1:H:153:ILE:HD11	1.85	0.59
1:G:89:PHE:CE1	1:G:145:VAL:HG11	2.40	0.56
1:H:98:LYS:CE	1:H:153:ILE:HD11	2.37	0.55
1:A:373:VAL:HG12	1:A:377:LEU:HD12	1.92	0.52
1:H:82:ILE:HG21	1:H:142:PHE:CD1	2.43	0.52
1:A:398:LEU:HG	1:A:399:LEU:HD23	1.92	0.51
1:H:98:LYS:HG2	1:H:153:ILE:CD1	2.42	0.50
1:H:94:MET:CE	1:H:145:VAL:HG12	2.43	0.49
1:H:83:PHE:HE2	1:H:142:PHE:HA	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:440:HIS:CE1	1:A:444:LEU:HD11	2.48	0.48
1:H:135:SER:HA	1:H:138:TRP:CE3	2.49	0.47
1:H:144:ARG:CB	1:H:203:MET:HE1	2.43	0.47
1:E:265:SER:HA	1:E:268:VAL:O	2.14	0.47
1:G:94:MET:SD	1:G:150:LEU:HD21	2.55	0.47
1:G:135:SER:HA	1:G:138:TRP:CE3	2.50	0.47
1:H:144:ARG:HA	1:H:203:MET:CE	2.44	0.46
1:E:19:GLU:HA	1:E:54:HIS:HB3	1.99	0.45
1:C:241:GLY:HA2	1:C:245:ARG:HB2	1.98	0.44
1:H:89:PHE:CE1	1:H:145:VAL:CG1	3.01	0.44
1:A:11:THR:HG21	1:A:398:LEU:HD13	2.00	0.43
1:B:39:ASP:HB3	1:D:313:TYR:CZ	2.54	0.43
1:B:135:SER:HA	1:B:138:TRP:CE3	2.53	0.43
1:D:241:GLY:HA2	1:D:245:ARG:HB2	2.01	0.43
1:B:396:TRP:HA	1:B:397:SER:HA	1.82	0.43
1:H:144:ARG:CA	1:H:203:MET:CE	2.96	0.43
1:F:435:ASN:O	1:F:439:THR:HG23	2.19	0.42
1:H:79:TRP:N	1:H:80:PRO:CD	2.82	0.42
1:F:241:GLY:HA2	1:F:245:ARG:HB2	2.01	0.42
1:G:139:PHE:CD2	1:G:196:SER:HB2	2.54	0.42
1:G:94:MET:HG2	1:G:149:GLU:HG3	2.00	0.42
1:A:135:SER:HA	1:A:138:TRP:CE3	2.54	0.42
1:B:3:PHE:HB2	1:B:440:HIS:CD2	2.54	0.42
1:E:79:TRP:HA	1:E:142:PHE:CE1	2.55	0.42
1:G:89:PHE:CE1	1:G:145:VAL:CG1	3.03	0.42
1:A:396:TRP:HA	1:A:397:SER:HA	1.86	0.42
1:G:204:LEU:HG	1:G:210:SER:HB2	2.02	0.41
1:H:82:ILE:HD11	1:H:145:VAL:CG1	2.38	0.41
1:H:82:ILE:CD1	1:H:145:VAL:HG11	2.37	0.41
1:C:293:PHE:CE2	1:C:330:PHE:HB2	2.55	0.41
1:H:396:TRP:HA	1:H:397:SER:HA	1.87	0.41
1:H:94:MET:HE3	1:H:145:VAL:HG12	2.02	0.41
1:C:135:SER:HA	1:C:138:TRP:CE3	2.56	0.41
1:G:140:LEU:HD11	1:G:199:LYS:HB3	2.04	0.40
1:C:396:TRP:HA	1:C:397:SER:HA	1.87	0.40
1:F:367:GLN:HA	1:F:370:ILE:HG13	2.04	0.40

All (1) 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)
1:B:207:GLU:OE2	1:H:88:GLN:NE2[2_1058]	1.81	0.39

### 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	442/471 (94%)	426 (96%)	16 (4%)	0	100	100
1	B	439/471 (93%)	426 (97%)	13 (3%)	0	100	100
1	C	441/471 (94%)	428 (97%)	13 (3%)	0	100	100
1	D	441/471 (94%)	426 (97%)	15 (3%)	0	100	100
1	E	436/471 (93%)	422 (97%)	14 (3%)	0	100	100
1	F	437/471 (93%)	423 (97%)	14 (3%)	0	100	100
1	G	439/471 (93%)	423 (96%)	16 (4%)	0	100	100
1	H	441/471 (94%)	424 (96%)	17 (4%)	0	100	100
All	All	3516/3768 (93%)	3398 (97%)	118 (3%)	0	100	100

There are no Ramachandran outliers to report.

#### 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	376/398 (94%)	373 (99%)	3 (1%)	81	87
1	B	375/398 (94%)	372 (99%)	3 (1%)	81	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	375/398 (94%)	374 (100%)	1 (0%)	92	95
1	D	375/398 (94%)	374 (100%)	1 (0%)	92	95
1	E	372/398 (94%)	367 (99%)	5 (1%)	69	76
1	F	373/398 (94%)	371 (100%)	2 (0%)	88	92
1	G	375/398 (94%)	370 (99%)	5 (1%)	69	76
1	H	375/398 (94%)	370 (99%)	5 (1%)	69	76
All	All	2996/3184 (94%)	2971 (99%)	25 (1%)	81	87

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASP
1	A	360	VAL
1	A	399	LEU
1	B	37	ASP
1	B	106	GLU
1	B	371	ASP
1	C	37	ASP
1	D	37	ASP
1	E	37	ASP
1	E	90	ASN
1	E	203	MET
1	E	204	LEU
1	E	275	HIS
1	F	209	ASN
1	F	439	THR
1	G	2	LYS
1	G	33	ASP
1	G	88	GLN
1	G	150	LEU
1	G	250	PRO
1	H	28	THR
1	H	37	ASP
1	H	98	LYS
1	H	204	LEU
1	H	371	ASP

Sometimes 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

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	$\# Z  > 2$
2	SO4	H	501	-	4,4,4	0.34	0	6,6,6	0.07	0
2	SO4	C	501	-	4,4,4	0.32	0	6,6,6	0.07	0
2	SO4	D	501	-	4,4,4	0.33	0	6,6,6	0.09	0
2	SO4	B	501	-	4,4,4	0.33	0	6,6,6	0.06	0
2	SO4	E	501	-	4,4,4	0.33	0	6,6,6	0.05	0

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.

## 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

### 6.1 Protein, DNA and RNA chains

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	444/471 (94%)	0.49	17 (3%) 40 40	26, 43, 63, 79	0
1	B	441/471 (93%)	0.37	7 (1%) 72 73	23, 40, 57, 77	0
1	C	443/471 (94%)	0.34	2 (0%) 91 91	27, 40, 56, 63	0
1	D	443/471 (94%)	0.25	3 (0%) 87 87	26, 36, 49, 57	0
1	E	438/471 (92%)	0.42	9 (2%) 63 65	27, 44, 66, 73	0
1	F	439/471 (93%)	0.29	3 (0%) 87 87	26, 42, 58, 64	0
1	G	441/471 (93%)	0.54	15 (3%) 45 45	32, 49, 73, 91	0
1	H	443/471 (94%)	0.63	18 (4%) 37 37	30, 49, 71, 82	0
All	All	3532/3768 (93%)	0.42	74 (2%) 63 65	23, 42, 66, 91	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	26	GLY	4.8
1	A	68	LEU	4.5
1	H	28	THR	4.4
1	G	7	PHE	4.1
1	A	360	VAL	4.0
1	H	88	GLN	3.9
1	A	419	PHE	3.8
1	F	228	ASP	3.6
1	H	82	ILE	3.4
1	D	5	PRO	3.4
1	G	6	ASN	3.3
1	A	442	ALA	3.3
1	A	443	ALA	3.3
1	G	89	PHE	3.1
1	A	356	ASP	3.0
1	B	55	TYR	3.0

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	H	142	PHE	3.0
1	E	154	VAL	3.0
1	H	95	ALA	2.9
1	A	364	ILE	2.8
1	A	9	PHE	2.8
1	H	89	PHE	2.7
1	G	441	LYS	2.7
1	H	97	TYR	2.7
1	A	3	PHE	2.7
1	E	141	ASP	2.6
1	B	126	GLU	2.6
1	F	315	ASP	2.6
1	G	5	PRO	2.6
1	A	361	ASP	2.6
1	A	2	LYS	2.6
1	G	3	PHE	2.6
1	G	394	TYR	2.6
1	A	399	LEU	2.5
1	A	445	PRO	2.5
1	G	420	ASP	2.5
1	E	438	GLN	2.5
1	H	141	ASP	2.5
1	H	135	SER	2.4
1	H	56	HIS	2.4
1	H	93	GLY	2.4
1	B	207	GLU	2.4
1	G	286	CYS	2.4
1	B	109	ILE	2.3
1	A	425	ILE	2.3
1	H	160	HIS	2.3
1	G	209	ASN	2.3
1	D	89	PHE	2.3
1	G	212	THR	2.3
1	E	37	ASP	2.2
1	E	385	ASP	2.2
1	A	420	ASP	2.2
1	H	59	GLU	2.2
1	E	200	ALA	2.2
1	H	204	LEU	2.2
1	B	108	GLY	2.2
1	G	144	ARG	2.1
1	B	9	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	359	LEU	2.1
1	C	361	ASP	2.1
1	A	106	GLU	2.1
1	C	441	LYS	2.1
1	E	122	MET	2.1
1	H	94	MET	2.1
1	E	136	VAL	2.1
1	F	436	VAL	2.1
1	B	441	LYS	2.1
1	G	119	ASP	2.1
1	H	124	ALA	2.1
1	G	145	VAL	2.1
1	E	3	PHE	2.1
1	D	444	LEU	2.1
1	H	200	ALA	2.1
1	G	380	VAL	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	SO4	B	501	5/5	0.90	0.18	67,68,68,69	0
2	SO4	H	501	5/5	0.91	0.25	59,60,60,61	0
2	SO4	E	501	5/5	0.92	0.14	69,69,70,71	0
2	SO4	C	501	5/5	0.94	0.15	56,56,57,58	0
2	SO4	D	501	5/5	0.95	0.19	53,53,54,55	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.