

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

#### Jul 31, 2023 – 08:58 PM EDT

PDB ID	:	5A3H
Title	:	2-DEOXY-2-FLURO-B-D-CELLOBIOSYL/ENZYME INTERMEDI-
		ATE COMPLEX OF THE ENDOGLUCANASE CEL5A FROM BACILLUS
		AGARADHEARANS AT 1.8 ANGSTROMS RESOLUTION
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		zie, L.; Withers, S.G.
Deposited on		
Resolution	:	1.82  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 (1)) were used in the production of this report:

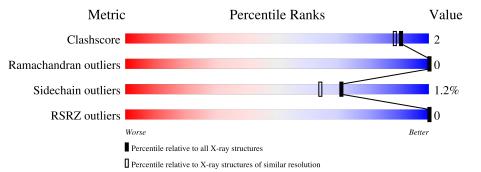
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as $541$ be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.34
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.34

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 1.82 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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 >=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	303	92%	6% ••			
2	В	2	50%	50%			



#### 5A3H

# 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2765 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 ENDOGLUCANASE.

Mol	Chain	Residues		Ate	oms			ZeroOcc	AltConf	Trace
1	А	300	Total 2394	C 1509	N 393	0 484	S 8	0	5	0

• Molecule 2 is an oligosaccharide called beta-D-glucopyranose-(1-4)-2-deoxy-2-fluoro-alpha-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	В	2	Total 23	C 12	F 1	O 10	0	1	0

• Molecule 3 is water.

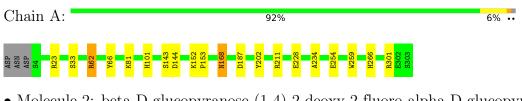
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	348	Total         O           348         348	0	0



# 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: ENDOGLUCANASE



• Molecule 2: beta-D-glucopyranose-(1-4)-2-deoxy-2-fluoro-alpha-D-glucopyranose

Chain B: 50% 50%



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	54.71Å 69.57Å 77.04Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	15.00 - 1.82	Depositor
Resolution (A)	14.90 - 1.80	EDS
% Data completeness	99.6 (15.00-1.82)	Depositor
(in resolution range)	97.2 (14.90-1.80)	EDS
R <sub>merge</sub>	0.06	Depositor
$R_{sym}$	0.06	Depositor
$< I/\sigma(I) > 1$	$2.97 (at 1.79 \text{\AA})$	Xtriage
Refinement program	REFMAC	Depositor
$R, R_{free}$	0.144 , $0.186$	Depositor
II, II, <i>free</i>	0.150 , (Not available)	DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor $(Å^2)$	11.7	Xtriage
Anisotropy	0.983	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.40 , $63.4$	EDS
L-test for twinning <sup>2</sup>	$ \langle L  \rangle = 0.49, \langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2765	wwPDB-VP
Average B, all atoms $(Å^2)$	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.30% 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.

# 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: G2F, BGC

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).

Mal	Chain	Bo	nd lengths	Bond angles		
Mol	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.64	2/2478~(0.1%)	1.17	10/3374~(0.3%)	

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	А	228	GLU	CD-OE2	10.72	1.37	1.25
1	А	228	GLU	CD-OE1	6.17	1.32	1.25

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	А	228	GLU	OE1-CD-OE2	-10.23	111.03	123.30
1	А	144	ASP	CB-CG-OD1	8.20	125.68	118.30
1	А	187	ASP	CB-CG-OD1	6.90	124.51	118.30
1	А	211	ARG	NE-CZ-NH1	6.45	123.53	120.30
1	А	62	ARG	NE-CZ-NH2	-6.16	117.22	120.30

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	2394	0	2203	8	0

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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 2.

The worst 5 of 8 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:152:LYS:HB3	1:A:153:PRO:HD3	1.92	0.51
1:A:254:GLU:HG3	3:A:1231:HOH:O	2.13	0.49
1:A:66:TYR:HA	1:A:101:HIS:HB3	1.98	0.45
1:A:33:SER:HA	1:A:62:ARG:HB3	2.01	0.42
1:A:202:TYR:CG	1:A:234:ALA:HA	2.55	0.42

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	6 Percentiles		
1	А	303/303~(100%)	293~(97%)	10 (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 side chain 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.



Chain Non-H H(model) H(added) Clashes Symm-Clashes Mol В  $\mathbf{2}$ 230 0 16 0 3 2 0 А 348 0 0 2765 All All 0 2219 8 0

The Analysed column shows the number of residues for which the side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	256/254 (101%)	253~(99%)	3 (1%)	71 64	

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	143	SER
1	А	168	ASN
1	А	266	HIS

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such side chains are listed below:

Mol	Chain	Res	Type
1	А	16	ASN
1	А	168	ASN

### 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)

3 monosaccharides 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	Mol Type Chain	Res	Link	Bond lengths			Bond angles			
10101		Ullalli	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
2	G2F	В	1[A]	-	11,11,12	0.91	0	$10,\!15,\!17$	1.07	0



	Mol T	Type Chain	Chain	ain Res	Res Link	Bo	Bond lengths			Bond angles		
		туре	Ullalli			Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
	2	G2F	В	1[B]	-	11,11,12	0.91	0	$10,\!15,\!17$	1.08	0	
	2	BGC	В	2	2	11,11,12	1.23	1 (9%)	$15,\!15,\!17$	1.81	6 (40%)	

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
2	G2F	В	1[A]	-	-	0/2/19/22	0/1/1/1
2	G2F	В	1[B]	-	-	0/2/19/22	0/1/1/1
2	BGC	В	2	2	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	В	2	BGC	O5-C1	-2.81	1.39	1.43

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms		$Observed(^{o})$	$Ideal(^{o})$
2	В	2	BGC	O4-C4-C3	-3.07	103.25	110.35
2	В	2	BGC	C2-C3-C4	-2.55	106.48	110.89
2	В	2	BGC	O5-C5-C4	-2.21	105.44	110.83
2	В	2	BGC	O3-C3-C2	-2.14	105.90	109.99
2	В	2	BGC	C1-O5-C5	2.08	115.02	112.19

There are no chirality outliers.

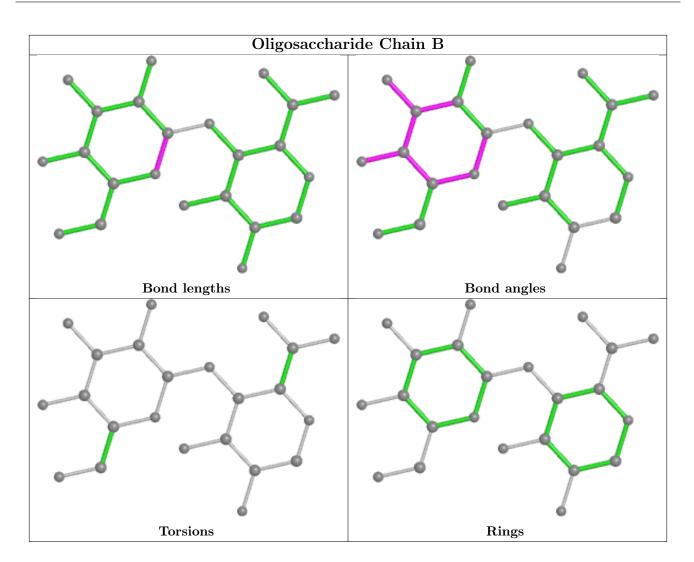
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





## 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	300/303~(99%)	-0.58	0 100 100	8, 12, 25, 41	1 (0%)

There are no RSRZ outliers to report.

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

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

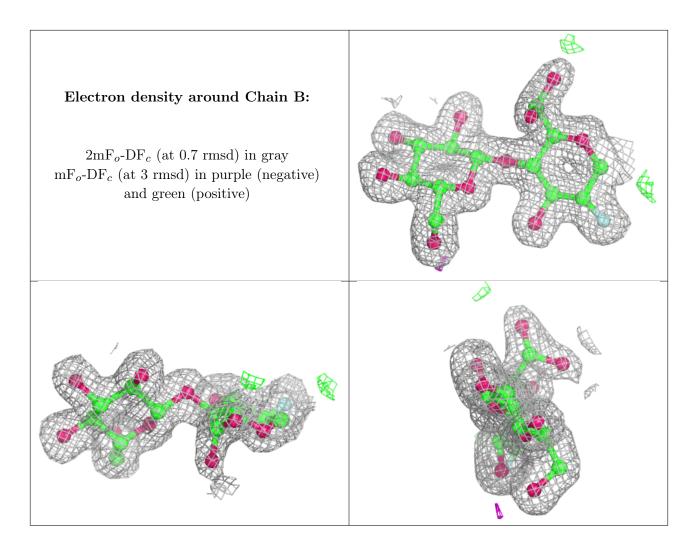
### 6.3 Carbohydrates (i)

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,  $95^{th}$  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	BGC	В	2	11/12	0.97	0.07	6, 9, 12, 14	0
2	G2F	В	1[B]	11/12	0.98	0.07	6, 9, 11, 11	1
2	G2F	В	1[A]	11/12	0.98	0.07	6,9,11,11	1

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.





## 6.4 Ligands (i)

There are no ligands in this entry.

## 6.5 Other polymers (i)

There are no such residues in this entry.

