

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

#### Oct 16, 2023 – 09:39 PM EDT

PDB ID : 1Z3T

Title: Structure of Phanerochaete chrysosporium cellobiohydrolase Cel7D (CBH58)

in complex with cellobiose

Authors: Ubhayasekera, W.; Stahlberg, J.; Mowbray, S.L.

Deposited on : 2005-03-14

Resolution : 1.70 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:

 $Mol Probity \quad : \quad 4.02b\text{--}467$ 

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

Xtriage (Phenix) : 1.13 EDS : 2.36

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)

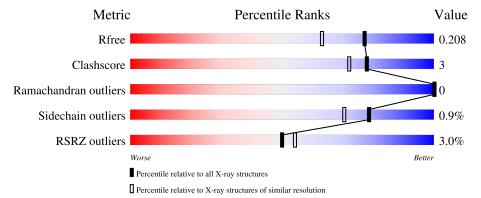
Validation Pipeline (wwPDB-VP) : 2.36

## 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 1.70 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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	A	431	89%	10%	•		
2	В	2	50%				



# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3340 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 cellulase.

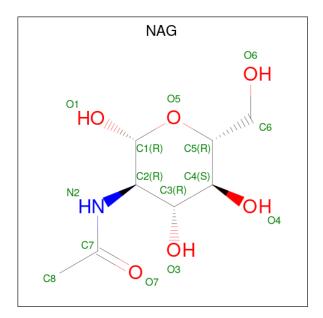
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	431	Total 3198	C 1962	N 541	O 667	S 28	0	0	0

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



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace	
2	В	2	Total 23	C 12	O 11	0	0	0

• Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 14 8 1 5	0	0

## • Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	105	Total O 105 105	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: cellulase

Chain A:

89%

10%

Molecule 2: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain B:

50%



# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	87.41Å 46.84Å 99.49Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 102.99° 90.00°	Depositor
Resolution (Å)	40.00 - 1.70	Depositor
Resolution (A)	39.36 - 1.70	EDS
% Data completeness	100.0 (40.00-1.70)	Depositor
(in resolution range)	99.6 (39.36-1.70)	EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	5.72 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
D D.	0.159 , 0.199	Depositor
$R, R_{free}$	0.176 , 0.208	DCC
$R_{free}$ test set	2541 reflections (5.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.6	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.36, 45.5	EDS
L-test for twinning <sup>2</sup>	$ < L > = 0.49, < L^2> = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3340	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.37% of the height of the origin peak. No significant pseudotranslation is detected.

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



<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: BGC, PCA, NAG

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	Bo	nd lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	1.25	$7/3256 \ (0.2\%)$	1.14	18/4429 (0.4%)	

#### All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	Observed(A)	$Ideal(\AA)$
1	A	376	SER	CB-OG	5.74	1.49	1.42
1	A	339	TRP	CG-CD1	-5.67	1.28	1.36
1	A	404	VAL	CB-CG2	-5.55	1.41	1.52
1	A	311	ALA	CA-CB	-5.34	1.41	1.52
1	A	114	TYR	CD1-CE1	-5.28	1.31	1.39
1	A	114	TYR	CE2-CZ	-5.13	1.31	1.38
1	A	168	TYR	CD2-CE2	-5.13	1.31	1.39

#### All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{\scriptscriptstyle o})$
1	A	147	ASP	CB-CG-OD2	11.65	128.79	118.30
1	A	381	ASP	CB-CG-OD2	8.09	125.58	118.30
1	A	63	ASP	CB-CG-OD2	7.40	124.96	118.30
1	A	251	ASP	CB-CG-OD1	-6.97	112.02	118.30
1	A	285	ASP	CB-CG-OD2	6.77	124.39	118.30
1	A	251	ASP	CB-CG-OD2	6.33	123.99	118.30
1	A	246	ASP	CB-CG-OD1	6.27	123.94	118.30
1	A	74	ASP	CB-CG-OD2	5.85	123.56	118.30
1	A	377	ASP	CB-CG-OD2	5.73	123.45	118.30
1	A	403	ASP	CB-CG-OD2	5.59	123.33	118.30
1	A	209	ASP	CB-CG-OD2	5.54	123.28	118.30
1	A	324	ASP	CB-CG-OD1	5.53	123.27	118.30
1	A	259	ASP	CB-CG-OD1	5.52	123.27	118.30
1	A	216	ASP	CB-CG-OD2	5.49	123.24	118.30

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Mol	Chain	Res	Type	Atoms	${f Z}$	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	336	ASP	CB-CG-OD2	5.30	123.07	118.30
1	A	149	ASP	CB-CG-OD2	5.14	122.93	118.30
1	A	422	ASP	CB-CG-OD1	5.08	122.87	118.30
1	A	366	ASP	CB-CG-OD1	5.03	122.82	118.30

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	A	3198	0	2974	18	0
2	В	23	0	21	0	0
3	A	14	0	13	0	0
4	A	105	0	0	0	0
All	All	3340	0	3008	18	0

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 3.

All (18) 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:284:ASN:HD21	1:A:291:THR:HB	1.37	0.90
1:A:284:ASN:ND2	1:A:291:THR:HB	1.91	0.85
1:A:407:GLN:HG2	1:A:408:VAL:HG13	1.72	0.71
1:A:317:ASP:H	1:A:329:GLN:HE22	1.42	0.65
1:A:317:ASP:H	1:A:329:GLN:NE2	1.96	0.64
1:A:227:THR:OG1	1:A:228:THR:N	2.27	0.60
1:A:404:VAL:HG23	1:A:408:VAL:CG2	2.35	0.57
1:A:18:GLN:HE21	1:A:28:LEU:HD21	1.74	0.53
1:A:25:CYS:O	1:A:26:LYS:HD3	2.13	0.48
1:A:404:VAL:HG23	1:A:408:VAL:HG22	1.94	0.48
1:A:222:PRO:HD2	1:A:250:CYS:O	2.16	0.45
1:A:284:ASN:HD21	1:A:291:THR:CB	2.17	0.44

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COHABABACA		DIEUIUU	DUIUE
0 0 1000100000			

Atom-1	Atom-2	Interatomic	$\operatorname{Clash}$	
Atom-1	Atom-2	${ m distance}({ m \AA})$	overlap (Å)	
1:A:383:ASP:OD1	1:A:384:PRO:HD2	2.18	0.44	
1:A:170:ASP:OD1	1:A:174:PRO:HD3	2.19	0.42	
1:A:250:CYS:HG	1:A:327:CYS:HA	1.84	0.42	
1:A:226:THR:HG23	1:A:343:LYS:HG3	2.02	0.41	
1:A:353:LEU:HD23	1:A:357:MET:SD	2.60	0.41	
1:A:250:CYS:SG	1:A:327:CYS:SG	3.20	0.40	

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	Percentiles	
1	A	429/431 (100%)	423 (99%)	6 (1%)	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	351/351 (100%)	348 (99%)	3 (1%)	78 70		

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



Mol	Chain	Res	Type
1	A	37	ASN
1	A	74	ASP
1	A	343	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	18	GLN
1	A	115	GLN
1	A	284	ASN
1	A	329	GLN
1	A	342	GLN

#### 5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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		Res Link		Bond lengths			Bond angles		
MOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	PCA	A	1	1	7,8,9	2.03	2 (28%)	9,10,12	2.13	4 (44%)

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
1	PCA	A	1	1	_	0/0/11/13	0/1/1/1



All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	Observed(A)	$Ideal(\AA)$
1	A	1	PCA	CA-N	4.38	1.51	1.46
1	A	1	PCA	CD-N	3.04	1.42	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Γype Atoms		$Observed(^o)$	$Ideal(^{o})$
1	A	1	PCA	CB-CA-C	-4.48	106.54	112.70
1	A	1	PCA	OE-CD-N	-2.42	119.22	124.86
1	A	1	PCA	CG-CD-N	2.29	114.31	108.39
1	A	1	PCA	CA-N-CD	-2.03	106.63	113.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates (i)

2 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 Type		Chain	Res	s Link	Bond lengths			Bond angles		
IVIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	BGC	В	1	2	12,12,12	0.81	0	17,17,17	0.77	0
2	BGC	В	2	2	11,11,12	1.03	0	15,15,17	1.42	2 (13%)

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	BGC	В	1	2	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	В	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
2	В	2	BGC	O2-C2-C1	3.18	115.65	109.15
2	В	2	BGC	C1-C2-C3	-2.43	106.68	109.67

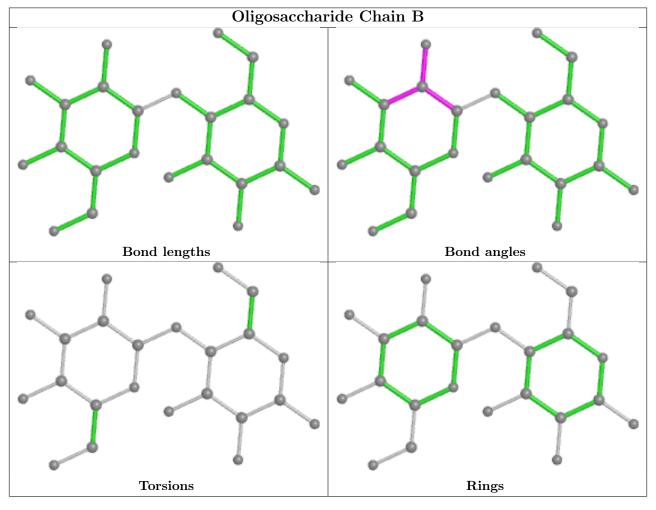
There are no chirality outliers.

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)

1 ligand is 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	Bo	ond leng	$ ag{ths}$	Bond angles		
	туре	Chain			Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	A	600	1	14,14,15	0.87	0	17,19,21	1.70	2 (11%)

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	NAG	A	600	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
3	A	600	NAG	O5-C1-C2	-4.66	103.93	111.29
3	A	600	NAG	O5-C5-C6	-2.48	103.31	107.20

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	600	NAG	O5-C5-C6-O6
3	A	600	NAG	C4-C5-C6-O6

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 (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	$\langle { m RSRZ} \rangle$	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	430/431 (99%)	0.14	13 (3%) 50 54	12, 24, 37, 46	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	431	SER	4.3
1	A	228	THR	4.1
1	A	236	ASP	4.0
1	A	383	ASP	3.1
1	A	45	SER	3.1
1	A	403	ASP	3.0
1	A	407	GLN	2.9
1	A	135	CYS	2.8
1	A	23	GLY	2.6
1	A	63	ASP	2.4
1	A	21	THR	2.3
1	A	25	CYS	2.2
1	A	313	ILE	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains (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	$\operatorname{Res}$	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{A}^2)$	Q<0.9
1	PCA	A	1	8/9	0.97	0.07	17,18,18,19	0

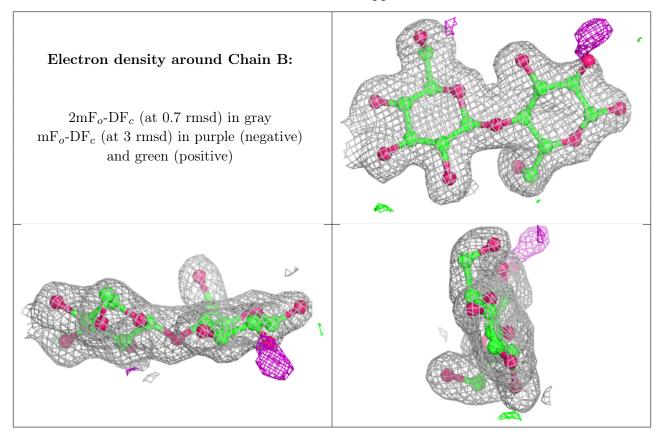


## 6.3 Carbohydrates (i)

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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	BGC	В	1	12/12	0.90	0.11	25,33,39,41	0
2	BGC	В	2	11/12	0.93	0.10	28,30,35,39	0

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)

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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
3	NAG	A	600	14/15	0.91	0.16	30,34,41,43	0



# 6.5 Other polymers (i)

There are no such residues in this entry.

