

Full wwPDB X-ray Structure Validation Report (i)

Oct 23, 2021 - 04:36 PM EDT

PDB ID : 1CEN

Title : CELLULASE (CELC) MUTANT WITH GLU 140 REPLACED BY GLN

COMPLEXED WITH CELLOHEXAOSE

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Deposited on : 1995-12-04

Resolution : 2.30 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 following versions of software and data (see references (i)) were used in the production of this report:

MolProbity: 4.02b-467

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

Xtriage (Phenix) : NOT EXECUTED EDS : NOT EXECUTED

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)

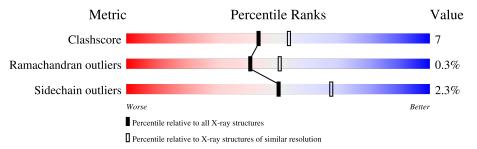
Validation Pipeline (wwPDB-VP) : 2.23.2

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 2.30 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	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain		
1	A	343	80%	17%	
2	В	2	100%		



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3048 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 CELC.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	334	Total	С	N	О	S	0	0	0
1	A	334	2804	1811	459	521	13	0	0	

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	140	GLN GLU e		engineered mutation	UNP P07985
A	156	CYS	TYR	conflict	UNP P07985

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



N	[ol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace	
	2	В	2	Total 23	C 12	O 11	0	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	221	Total O 221 221	0	0

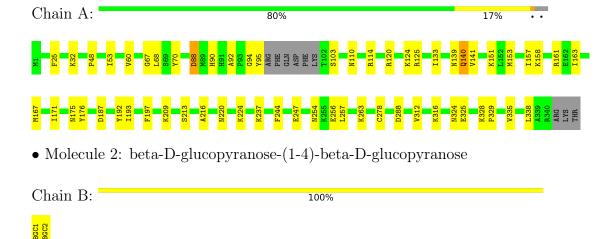


3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

• Molecule 1: CELLULASE CELC





4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source	
Space group	P 21 21 21	Depositor	
Cell constants	51.40Å 84.33Å 87.50Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	10.00 - 2.30	Depositor	
% Data completeness	(Not available) (10.00-2.30)	Depositor	
(in resolution range)	(10.00 2.50)		
R_{merge}	0.09	Depositor	
R_{sym}	(Not available)	Depositor	
Refinement program	X-PLOR	Depositor	
R, R_{free}	0.178 , 0.246	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	3048	wwPDB-VP	
Average B, all atoms (Å ²)	22.0	wwPDB-VP	



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

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).

\ ₁	Mol	Chain	Bond	$\mathbf{lengths}$	Bond angles		
ľ			RMSZ	# Z > 5	RMSZ	# Z > 5	
	1	A	0.45	0/2879	0.65	0/3888	

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	2804	0	2699	37	0
2	В	23	0	21	0	0
3	A	221	0	0	8	0
All	All	3048	0	2720	37	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 7.

All (37) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} \operatorname{Clash} \ \operatorname{overlap}\ (\mathring{\mathbf{A}}) \end{aligned}$
1:A:171:ILE:HD12	1:A:193:ILE:HG23	1.75	0.69

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Atom-1	Atom-2	Interatomic	Clash
		distance (Å)	overlap (Å)
1:A:167:MET:SD	3:A:617:HOH:O	2.50	0.68
1:A:161:ARG:NH2	1:A:192:TYR:HB2	2.10	0.67
1:A:120:ARG:HD3	3:A:436:HOH:O	1.98	0.62
1:A:175:ASN:HB3	3:A:598:HOH:O	2.01	0.60
1:A:153:MET:O	1:A:157:ILE:HG12	2.04	0.57
1:A:324:ASN:HD22	1:A:328:LYS:HB2	1.70	0.55
1:A:175:ASN:HB3	3:A:611:HOH:O	2.08	0.53
1:A:140:GLN:HB3	1:A:176:TYR:HA	1.90	0.52
1:A:161:ARG:HH22	1:A:192:TYR:HB2	1.74	0.52
1:A:312:VAL:HG11	1:A:338:LEU:HD13	1.92	0.51
1:A:197:PHE:O	1:A:278:CYS:HA	2.11	0.50
1:A:90:HIS:HA	1:A:139:ASN:HB3	1.93	0.49
1:A:95:TYR:OH	1:A:141:VAL:HA	2.11	0.49
1:A:167:MET:CE	3:A:617:HOH:O	2.62	0.48
1:A:53:ILE:O	1:A:67:GLY:HA3	2.14	0.48
1:A:224:LYS:HD2	1:A:288:ASP:OD1	2.14	0.47
1:A:94:GLY:O	1:A:103:SER:HA	2.15	0.46
1:A:124:LYS:HG3	1:A:163:ILE:HD13	1.97	0.46
1:A:110:ASN:O	1:A:114:ARG:HG3	2.15	0.45
1:A:213:SER:HB3	1:A:216:ALA:HB3	1.99	0.45
1:A:244:PHE:O	1:A:247:GLU:HG2	2.17	0.45
1:A:151:LYS:HE3	3:A:473:HOH:O	2.17	0.44
1:A:133:ILE:HG23	1:A:167:MET:HE1	1.98	0.44
1:A:329:PRO:HB3	1:A:335:VAL:HG21	1.99	0.43
1:A:171:ILE:HD12	1:A:193:ILE:CG2	2.45	0.43
1:A:254:ASN:O	1:A:257:LEU:HB3	2.18	0.43
1:A:209:LYS:HE2	1:A:220:ASN:OD1	2.18	0.43
1:A:237:LYS:HG2	3:A:599:HOH:O	2.18	0.42
1:A:60:VAL:HG11	1:A:114:ARG:HG2	2.02	0.42
1:A:256:GLU:HG3	3:A:453:HOH:O	2.19	0.42
1:A:48:PRO:HA	1:A:88:ASP:O	2.21	0.41
1:A:140:GLN:CB	1:A:176:TYR:HA	2.51	0.41
1:A:68:LEU:HD13	1:A:125:ARG:NH1	2.36	0.40
1:A:171:ILE:CD1	1:A:193:ILE:HD12	2.51	0.40
1:A:25:PHE:HB3	1:A:70:TYR:CZ	2.57	0.40
1:A:92:ALA:HB3	1:A:95:TYR:HB3	2.03	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.

Mo	ol Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	330/343 (96%)	318 (96%)	11 (3%)	1 (0%)	41 50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	GLN

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 O		Outliers	Percentiles
1	A	302/311 (97%)	295 (98%)	7 (2%)	50 67

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LYS
1	A	88	ASP
1	A	158	LYS
1	A	187	ASP
1	A	263	LYS
1	A	316	LYS
1	A	325	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	16	GLN
1	A	112	GLN
1	A	140	GLN
1	A	201	ASN
1	A	324	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)

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	Chain	Chain	Chain	Res	Link	Bo	ond leng	$ ag{ths}$	В	ond ang	cles
WIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2		
2	BGC	В	1	2	12,12,12	1.62	1 (8%)	17,17,17	2.06	8 (47%)		
2	BGC	В	2	2	11,11,12	1.32	1 (9%)	15,15,17	1.28	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
2	BGC	В	2	2	-	0/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
2	В	1	BGC	C1-C2	4.77	1.63	1.52
2	В	2	BGC	C4-C3	3.27	1.60	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
2	В	1	BGC	O2-C2-C1	3.86	118.11	109.16
2	В	1	BGC	C4-C3-C2	-3.77	104.23	110.82
2	В	1	BGC	O2-C2-C3	-2.57	104.41	110.35
2	В	2	BGC	O5-C1-C2	2.45	114.55	110.77
2	В	2	BGC	C1-O5-C5	2.39	115.43	112.19
2	В	1	BGC	O3-C3-C4	2.36	115.80	110.35
2	В	1	BGC	C3-C4-C5	-2.20	106.32	110.24
2	В	1	BGC	C6-C5-C4	2.17	118.08	113.00
2	В	1	BGC	O5-C1-C2	2.12	114.07	110.28
2	В	1	BGC	O5-C5-C4	-2.02	106.03	109.69

There are no chirality outliers.

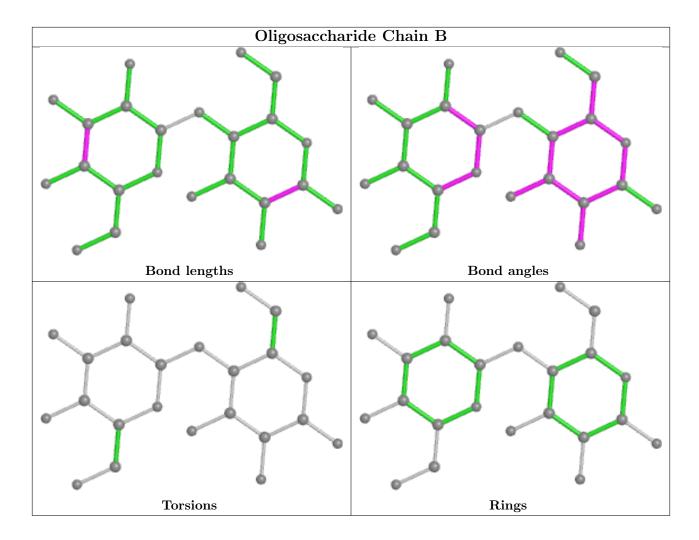
There are no torsion outliers.

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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

