



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2018 – 12:21 am GMT

PDB ID : 1EO7  
Title : BACILLUS CIRCULANS STRAIN 251 CYCLODEXTRIN GLYCOSYL-TRANSFERASE IN COMPLEX WITH MALTOHEXAOSE  
Authors : Uitdehaag, J.C.M.; Dijkstra, B.W.  
Deposited on : 2000-03-22  
Resolution : 2.48 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix)	:	1.13
EDS	:	trunk30967
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30967

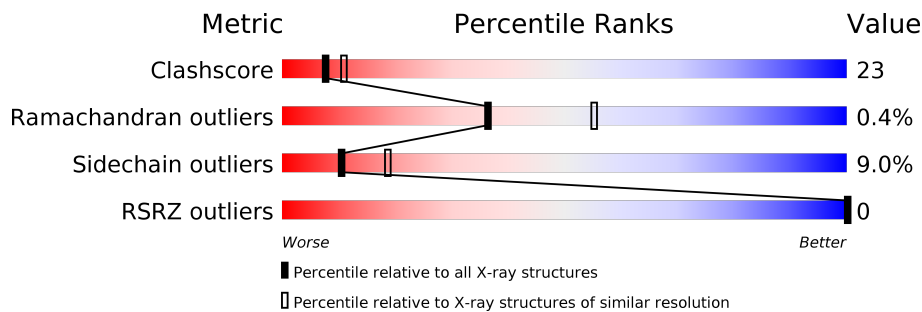
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.48 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))
Clashscore	122126	5860 (2.50-2.46)
Ramachandran outliers	120053	5763 (2.50-2.46)
Sidechain outliers	120020	5765 (2.50-2.46)
RSRZ outliers	108989	5026 (2.50-2.46)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	686	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	A	691	X	-	-	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5585 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.

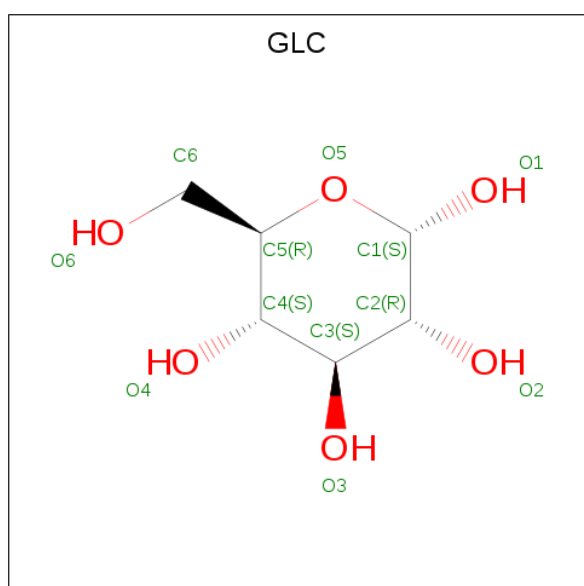
- Molecule 1 is a protein called PROTEIN (CYCLODEXTRIN GLYCOSYLTRANSFERASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	686	Total	C	N	O	S	0	0	0
			5257	3318	900	1024	15			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	229	ALA	ASP	ENGINEERED	UNP P43379
A	257	ALA	GLU	ENGINEERED	UNP P43379
A	400	SER	CYS	CONFLICT	UNP P43379

- Molecule 2 is ALPHA-D-GLUCOSE (three-letter code: GLC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			11	6	5		
2	A	1	Total	C	O	0	0
			11	6	5		

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 12 6 6	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 12 6 6	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 12 6 6	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 12 6 6	0	0
2	A	1	Total C O 11 6 5	0	0
2	A	1	Total C O 12 6 6	0	0

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	3	Total Ca 3 3	0	0

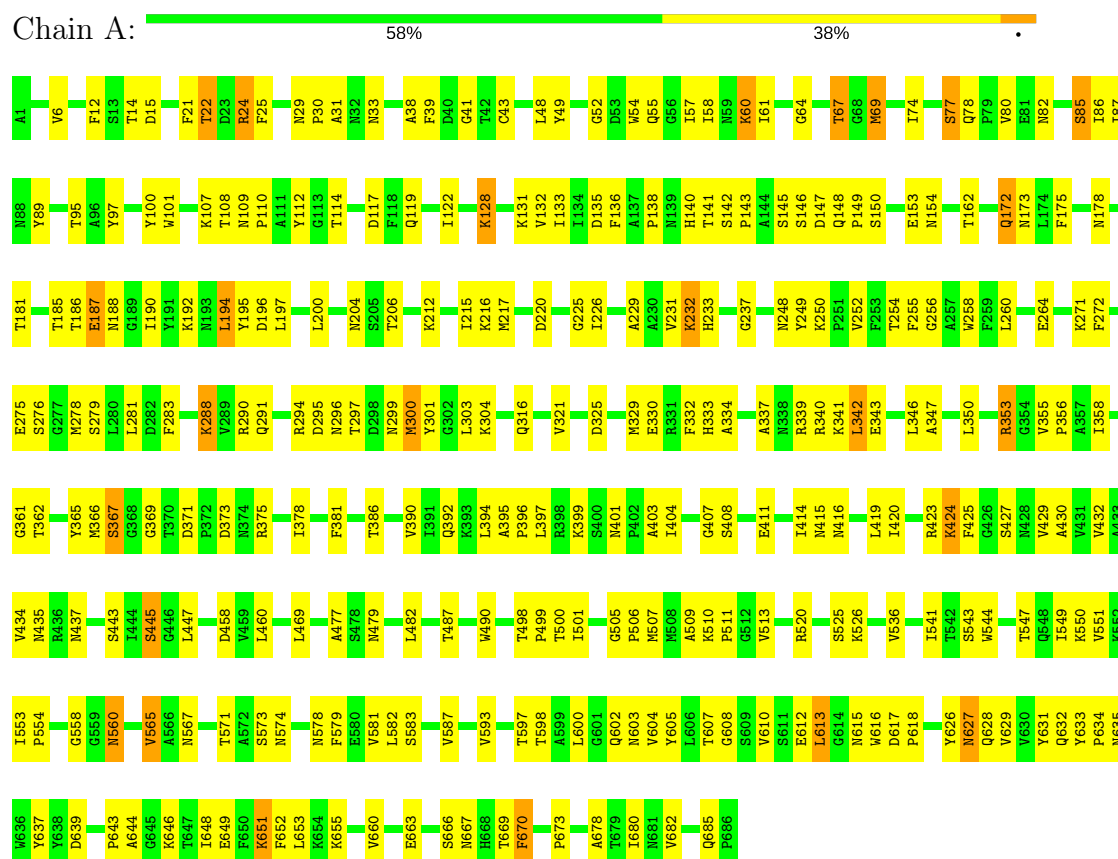
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	145	Total O 145 145	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: PROTEIN (CYCLODEXTRIN GLYCOSYLTRANSFERASE)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.65Å 109.02Å 64.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.48 78.00 – 2.48	Depositor EDS
% Data completeness (in resolution range)	91.0 (7.00-2.48) 88.3 (78.00-2.48)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.69 (at 2.48Å)	Xtriage
Refinement program	TNT	Depositor
R, $R_{free}$	0.228 , 0.296 0.211 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.0	Xtriage
Anisotropy	0.312	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 62.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5585	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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

### 5.1 Standard geometry

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

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	0.36	0/5387	0.50	0/7343

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

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	5257	0	5025	244	0
2	A	180	0	156	7	0
3	A	3	0	0	0	0
4	A	145	0	0	11	0
All	All	5585	0	5181	247	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 23.

The worst 5 of 247 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:330:GLU:HG2	1:A:369:GLY:HA2	1.34	1.10
1:A:38:ALA:HB2	1:A:86:ILE:HD11	1.39	1.04
1:A:536:VAL:HG21	1:A:554:PRO:HG3	1.46	0.93
1:A:300:MET:HG3	1:A:419:LEU:HB2	1.50	0.92
1:A:231:VAL:HG11	1:A:256:GLY:HA3	1.56	0.87

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	684/686 (100%)	637 (93%)	44 (6%)	3 (0%)	36 55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	629	VAL
1	A	22	THR
1	A	627	ASN

### 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	554/554 (100%)	504 (91%)	50 (9%)	10 19



5 of 50 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	217	MET
1	A	342	LEU
1	A	660	VAL
1	A	232	LYS
1	A	288	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	269	ASN
1	A	479	ASN
1	A	578	ASN
1	A	263	ASN
1	A	560	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](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 3 are monoatomic - leaving 16 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
2	GLC	A	691	2	11,11,12	0.60	0	15,15,17	3.11	5 (33%)
2	GLC	A	692	2	11,11,12	0.47	0	15,15,17	0.92	1 (6%)
2	GLC	A	693	2	11,11,12	0.45	0	15,15,17	0.79	0
2	GLC	A	694	2	11,11,12	0.36	0	15,15,17	0.87	1 (6%)
2	GLC	A	695	2	11,11,12	0.36	0	15,15,17	0.69	0
2	GLC	A	696	2	12,12,12	0.32	0	17,17,17	0.65	0
2	GLC	A	698	2	11,11,12	0.34	0	15,15,17	0.74	0
2	GLC	A	699	2	11,11,12	0.36	0	15,15,17	0.67	0
2	GLC	A	700	2	12,12,12	0.31	0	17,17,17	0.83	0
2	GLC	A	702	2	11,11,12	0.35	0	15,15,17	0.91	0
2	GLC	A	703	2	11,11,12	0.36	0	15,15,17	0.65	0
2	GLC	A	704	2	11,11,12	0.36	0	15,15,17	0.66	0
2	GLC	A	705	2	12,12,12	0.31	0	17,17,17	0.87	0
2	GLC	A	707	2	11,11,12	0.32	0	15,15,17	0.64	0
2	GLC	A	708	2	11,11,12	0.35	0	15,15,17	0.70	0
2	GLC	A	709	2	12,12,12	0.31	0	17,17,17	0.95	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
2	GLC	A	691	2	1/1/4/5	0/2/18/22	0/1/1/1
2	GLC	A	692	2	-	0/2/18/22	0/1/1/1
2	GLC	A	693	2	-	0/2/18/22	0/1/1/1
2	GLC	A	694	2	-	0/2/18/22	0/1/1/1
2	GLC	A	695	2	-	0/2/18/22	0/1/1/1
2	GLC	A	696	2	-	0/2/22/22	0/1/1/1
2	GLC	A	698	2	-	0/2/18/22	0/1/1/1
2	GLC	A	699	2	-	0/2/18/22	0/1/1/1
2	GLC	A	700	2	-	0/2/22/22	0/1/1/1
2	GLC	A	702	2	-	0/2/18/22	0/1/1/1
2	GLC	A	703	2	-	0/2/18/22	0/1/1/1
2	GLC	A	704	2	-	0/2/18/22	0/1/1/1
2	GLC	A	705	2	-	0/2/22/22	0/1/1/1
2	GLC	A	707	2	-	0/2/18/22	0/1/1/1
2	GLC	A	708	2	-	0/2/18/22	0/1/1/1
2	GLC	A	709	2	-	0/2/22/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	691	GLC	C4-C3-C2	-7.16	101.17	110.61
2	A	691	GLC	O2-C2-C1	-5.96	95.16	109.14
2	A	691	GLC	O5-C5-C4	-4.26	103.02	109.97
2	A	691	GLC	O3-C3-C2	-2.21	105.88	110.19
2	A	709	GLC	O5-C5-C4	2.02	113.39	109.69

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	691	GLC	C1

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	691	GLC	3	0
2	A	692	GLC	1	0
2	A	693	GLC	1	0
2	A	694	GLC	1	0
2	A	695	GLC	2	0
2	A	696	GLC	2	0

## 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<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	686/686 (100%)	-0.19	0 100 100	29, 41, 54, 72	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](#)

There are no carbohydrates 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	GLC	A	691	11/12	0.76	0.28	12,14,20,21	0
2	GLC	A	698	11/12	0.81	0.24	38,38,42,45	0
2	GLC	A	709	12/12	0.82	0.17	51,53,62,65	0
2	GLC	A	702	11/12	0.84	0.20	65,66,71,74	0
2	GLC	A	705	12/12	0.84	0.25	62,64,72,100	0
2	GLC	A	693	11/12	0.87	0.13	47,49,54,57	0
2	GLC	A	703	11/12	0.90	0.16	61,61,66,69	0
2	GLC	A	700	12/12	0.91	0.13	39,40,46,49	0
2	GLC	A	694	11/12	0.91	0.14	49,50,55,59	0

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GLC	A	699	11/12	0.92	0.16	41,43,48,51	0
2	GLC	A	708	11/12	0.92	0.15	47,49,54,56	0
2	GLC	A	695	11/12	0.92	0.14	49,51,56,59	0
2	GLC	A	704	11/12	0.93	0.15	59,61,66,69	0
2	GLC	A	696	12/12	0.93	0.14	50,51,59,61	0
2	GLC	A	707	11/12	0.94	0.15	46,46,51,54	0
3	CA	A	688	1/1	0.95	0.09	57,57,57,57	0
2	GLC	A	692	11/12	0.96	0.11	33,46,51,54	0
3	CA	A	690	1/1	0.96	0.14	52,52,52,52	0
3	CA	A	689	1/1	0.96	0.12	39,39,39,39	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.