

wwPDB X-ray Structure Validation Summary Report (i)

Oct 30, 2023 - 07:10 PM JST

:	4XWM
:	Complex structure of catalytic domain of Clostridium Cellulovorans Exgs and
	Cellobiose
:	Liaw, YC.
:	2015-01-29
:	1.70 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:

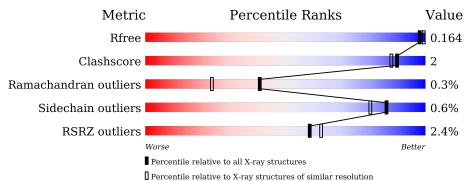
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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: $X\text{-}RAY \, DIFFRACTION$

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	$\begin{array}{l} \textbf{Whole archive} \\ \textbf{(\#Entries)} \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
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	А	681	2% 	• 7%
2	В	2	100%	
3	С	2	100%	
3	D	2	50% 50%	



4XWM

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5828 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 Exoglucanase S.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	630	Total 5152	C 3310	N 828	O 996	S 18	0	38	0

Chain	Residue	Modelled	Actual	Comment	Reference
А	-43	MET	-	expression tag	UNP O65986
А	-42	GLY	-	expression tag	UNP O65986
А	-41	SER	-	expression tag	UNP O65986
А	-40	SER	-	expression tag	UNP O65986
А	-39	HIS	-	expression tag	UNP O65986
А	-38	HIS	-	expression tag	UNP O65986
А	-37	HIS	-	expression tag	UNP O65986
A	-36	HIS	-	expression tag	UNP O65986
А	-35	HIS	-	expression tag	UNP O65986
A	-34	HIS	-	expression tag	UNP O65986
A	-33	SER	-	expression tag	UNP O65986
А	-32	SER	-	expression tag	UNP O65986
A	-31	GLY	-	expression tag	UNP O65986
А	-30	LEU	-	expression tag	UNP O65986
A	-29	VAL	-	expression tag	UNP O65986
А	-28	PRO	-	expression tag	UNP O65986
А	-27	ARG	-	expression tag	UNP O65986
A	-26	GLY	-	expression tag	UNP O65986
А	-25	SER	-	expression tag	UNP O65986
A	-24	HIS	-	expression tag	UNP O65986
А	-23	MET	-	expression tag	UNP O65986
A	-22	ALA	-	expression tag	UNP O65986
А	-21	SER	-	expression tag	UNP O65986
А	-20	MET	-	expression tag	UNP O65986
А	-19	THR	-	expression tag	UNP O65986
А	-18	GLY	-	expression tag	UNP O65986
А	-17	GLY	_	expression tag	UNP O65986

There are 38 discrepancies between the modelled and reference sequences:

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Chain	Chain Residue Modelled Actual Comment							
Unam	nesique	Modelled	Actual	Comment	Reference			
A	-16	GLN	-	expression tag	UNP O65986			
А	-15	GLN	-	expression tag	UNP O65986			
А	-14	MET	-	expression tag	UNP O65986			
A	-13	GLY	-	expression tag	UNP O65986			
А	-12	ARG	-	expression tag	UNP O65986			
A	-11	ILE	-	expression tag	UNP O65986			
A	-10	GLU	-	expression tag	UNP O65986			
А	-9	GLY	-	expression tag	UNP O65986			
А	-8	ARG	-	expression tag	UNP O65986			
А	-7	GLU	-	expression tag	UNP O65986			
А	-6	PHE	-	expression tag	UNP O65986			

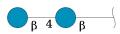
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• Molecule 2 is an oligosaccharide called beta-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	At	\mathbf{oms}		ZeroOcc	AltConf	Trace
2	В	2	Total 23	C 12	0 11	0	0	0

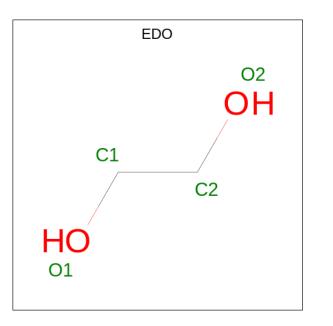
• Molecule 3 is an oligosaccharide called beta-D-glucopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
3	С	2	Total C O 23 12 11	0	0	0
3	D	2	Total C O 23 12 11	0	0	0

• Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	А	1	Total 4	${ m C} 2$	O 2	0	0

• Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	2	Total Ca 2 2	0	0

• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	597	Total O 601 601	0	4



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.

Chain A:	89%	• 7%
MET GLY SER SER SER HIS HIS HIS HIS SER SER	GLY VAL VAL PRO ARG GLY GLY MET HIS MET THR MET GLN GLN GLN GLN GLN GLN ARG GLU GLN ARG GLU GLN PHE FRO FRO FRO FRO FRO	VAL V-1 P0 N1 E39 A40 P41 H45 P111 B3 M115 D154
1166 W241 1253 W290 P332 V394	K432 W468 W469 Y469 Y469 K512 D508 D508 D508 K512 K512 K512 K512 K564 K512 K564 K512 K564 K598 K598 K598 K598 K598 K598 K501 K501	Ceo2 Veo3 PRO PRO PRO CLY LVS CLY CLY CLY CLY CLY
• Molecule 2: be	eta-D-glucopyranose-(1-4)-alpha-D-gluco	pyranose
Chain B:	100%	
GLC1 BGC2		
• Molecule 3: be	eta-D-glucopyranose-(1-4)-beta-D-glucop	oyranose
Chain C:	100%	
BGC1 BGC2		
• Molecule 3: be	eta-D-glucopyranose-(1-4)-beta-D-glucop	byranose
Chain D:	50%	50%
BGC1 BGC2		

 \bullet Molecule 1: Exoglucan ase S



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants	109.14Å 109.14Å 182.15Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.50 - 1.70	Depositor
Resolution (A)	27.51 - 1.70	EDS
% Data completeness	99.5 (27.50-1.70)	Depositor
(in resolution range)	97.8 (27.51-1.70)	EDS
R _{merge}	0.05	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$6.14 (at 1.71 \text{\AA})$	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
D D.	0.146 , 0.163	Depositor
R, R_{free}	0.148 , 0.164	DCC
R_{free} test set	2000 reflections (1.67%)	wwPDB-VP
Wilson B-factor $(Å^2)$	11.4	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.39 , 48.0	EDS
L-test for twinning ²	$< L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5828	wwPDB-VP
Average B, all atoms $(Å^2)$	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.53% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities 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: EDO, GLC, BGC, CA

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		lengths		angles
Mol Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.40	0/5424	0.57	0/7386

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	45	HIS	Peptide
1	А	599	ASN	Sidechain

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	5152	0	4860	18	0

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Mol	Chain	-	H(model)	H(added)	Clashes	Symm-Clashes
2	В	23	0	21	0	0
3	С	23	0	21	0	0
3	D	23	0	21	0	0
4	А	4	0	6	0	0
5	А	2	0	0	0	0
6	А	601	0	0	1	0
All	All	5828	0	4929	18	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 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:506:SER:O	1:A:512[B]:LYS:HE3	2.00	0.62
1:A:253[B]:THR:CG2	1:A:332:PRO:HA	2.34	0.58
1:A:40:ALA:HB3	1:A:41:PRO:HD3	1.88	0.56
1:A:564:LYS:HE3	6:A:938:HOH:O	2.06	0.55
1:A:508:ASP:O	1:A:512[B]:LYS:HG2	2.06	0.55

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	А	666/681~(98%)	652 (98%)	12 (2%)	2~(0%)	41 24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type		
1	А	40	ALA		
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Mol	Chain	Res	Type
1	А	39	GLU

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.

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	А	542/550~(98%)	538~(99%)	4 (1%)	84 77

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	83[A]	ILE
1	А	83[B]	ILE
1	А	584	TYR
1	А	599	ASN

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

Mol	Chain	Res	Type
1	А	326	ASN
1	А	342	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)

6 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	Turne	Chain	Res	Link	Bo	ond leng	ths	Bond angles		
	Type	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
2	GLC	В	1	2	12,12,12	0.49	0	17,17,17	0.78	0
2	BGC	В	2	2	11,11,12	0.43	0	$15,\!15,\!17$	0.63	0
3	BGC	С	1	3	12,12,12	0.42	0	17,17,17	0.78	0
3	BGC	С	2	3	11,11,12	0.28	0	$15,\!15,\!17$	1.00	0
3	BGC	D	1	3	12,12,12	0.48	0	$17,\!17,\!17$	0.73	0
3	BGC	D	2	3	11,11,12	0.25	0	$15,\!15,\!17$	1.34	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	GLC	В	1	2	-	0/2/22/22	0/1/1/1
2	BGC	В	2	2	-	0/2/19/22	0/1/1/1
3	BGC	С	1	3	-	0/2/22/22	0/1/1/1
3	BGC	С	2	3	-	0/2/19/22	0/1/1/1
3	BGC	D	1	3	-	1/2/22/22	0/1/1/1
3	BGC	D	2	3	-	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	$Observed(^{o})$	$Ideal(^{o})$
3	D	2	BGC	C1-O5-C5	3.20	116.53	112.19
3	D	2	BGC	C2-C3-C4	-2.26	106.99	110.89

There are no chirality outliers.

All (1) torsion outliers are listed below:

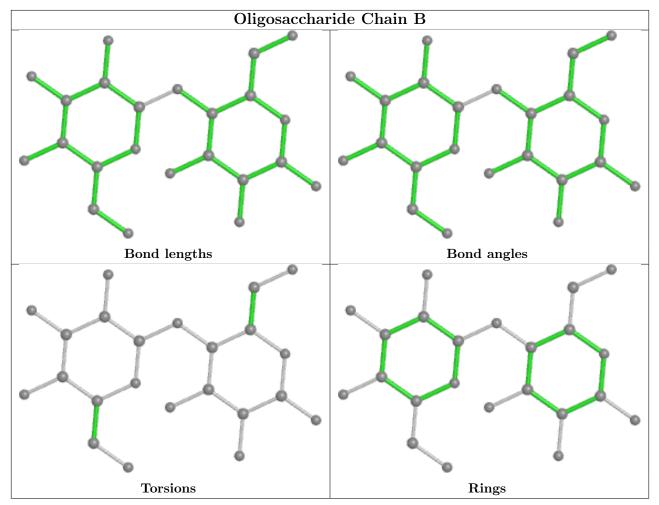
Mol	Chain	Res	Type	Atoms
3	D	1	BGC	O5-C5-C6-O6



There are no ring outliers.

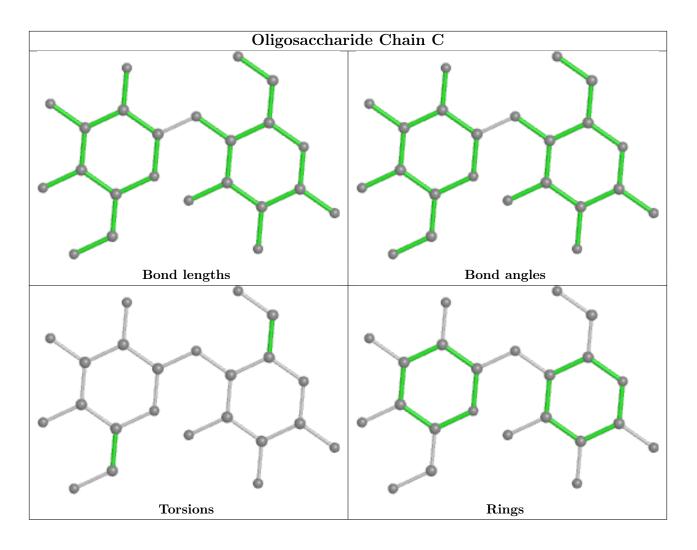
No monomer is involved in short contacts.

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



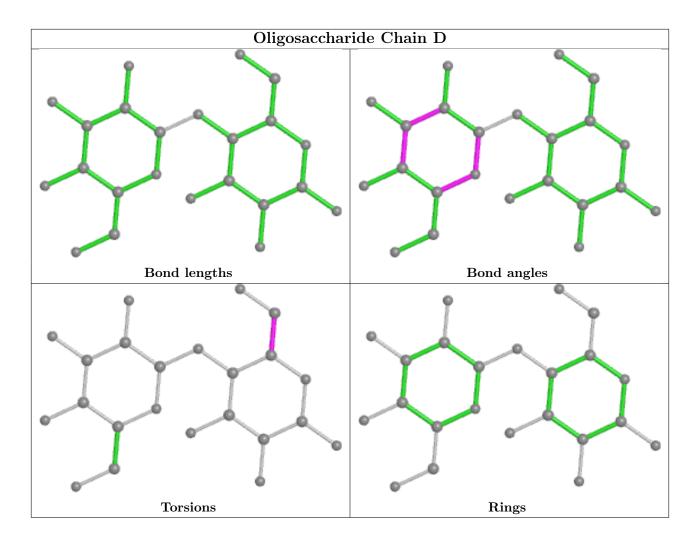












5.6 Ligand geometry (i)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	Chain	Res	Link	Bond lengths			Bond angles			
	туре	Unam	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	А	702	-	3,3,3	0.45	0	$2,\!2,\!2$	0.30	0

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
4	EDO	A	702	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

	Mol	Chain	Res	Type	Atoms
ſ	4	А	702	EDO	O1-C1-C2-O2

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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	Q<0.9
1	А	630/681~(92%)	-0.30	15 (2%) 59 63	5, 11, 26, 63	0

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	600	GLY	9.4
1	А	-1	VAL	8.3
1	А	598	ALA	6.4
1	А	599	ASN	6.4
1	А	601	THR	6.2

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)

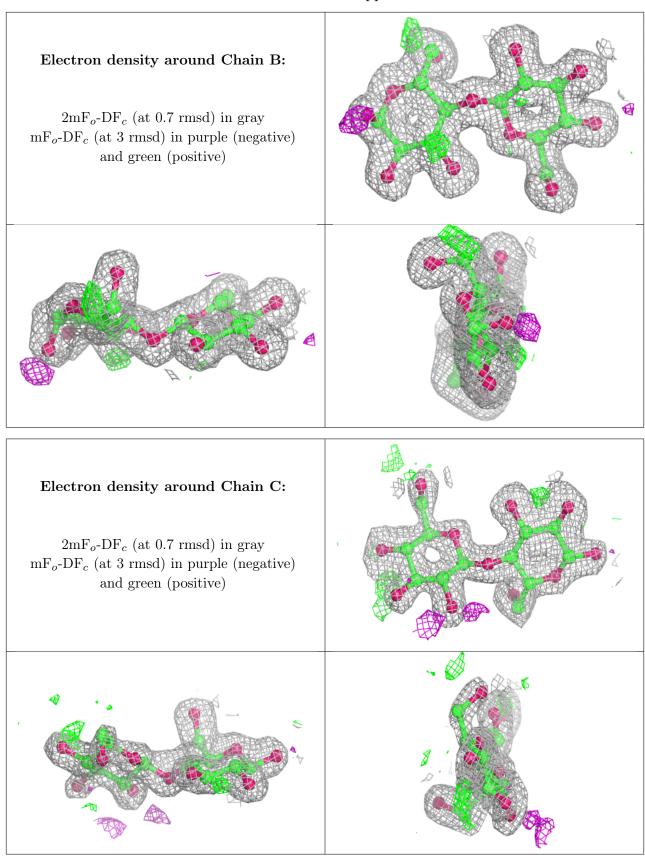
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(Å^2)$	Q < 0.9
3	BGC	D	2	11/12	0.59	0.29	54,59,66,67	0
3	BGC	D	1	12/12	0.82	0.19	25,47,52,57	0
3	BGC	С	2	11/12	0.88	0.15	16,21,27,30	11
3	BGC	С	1	12/12	0.94	0.10	16,18,20,21	12
2	GLC	В	1	12/12	0.96	0.07	11,16,21,29	0
2	BGC	В	2	11/12	0.98	0.05	8,10,12,14	0

The following is a graphical depiction of the model fit to experimental electron density for oligosac-

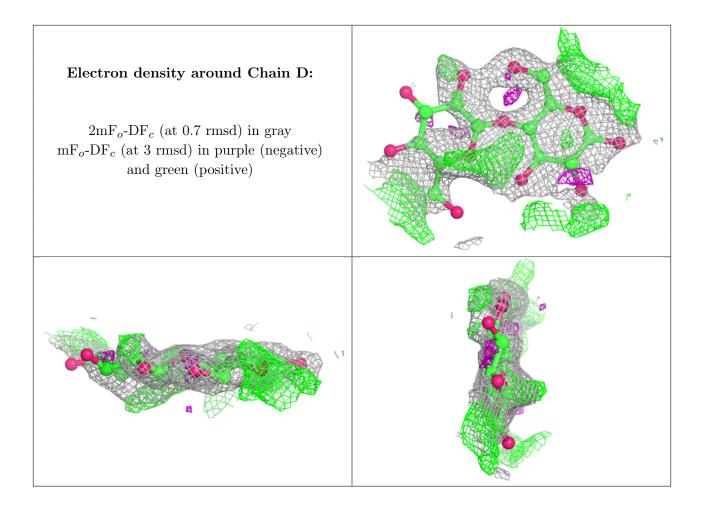






charide. Each fit is shown from different orientation to approximate a three-dimensional view.





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	B-factors(Å ²)	Q<0.9
4	EDO	А	702	4/4	0.71	0.22	44,50,56,58	0
5	CA	А	705	1/1	0.97	0.16	48,48,48,48	0
5	CA	А	706	1/1	0.99	0.16	18,18,18,18	0

6.5 Other polymers (i)

There are no such residues in this entry.

