



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 15, 2020 – 10:34 AM BST

PDB ID : 7C6N
Title : Crystal structure of beta-glycosides-binding protein (W177X) of ABC transporter in a closed state bound to cellotetraose (Form II)
Authors : Kanaujia, S.P.; Chandravanshi, M.; Samanta, R.
Deposited on : 2020-05-21
Resolution : 2.05 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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 : 2.14.4.dev1
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) : 2.14.4.dev1

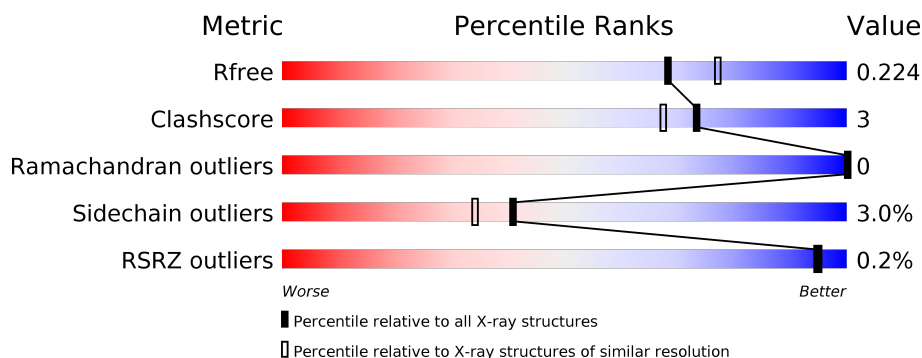
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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))
R_{free}	130704	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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 ≥ 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 $\leq 5\%$. 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	422	<div> <div>88%</div> <div>10% ..</div> </div>
1	B	422	<div> <div>88%</div> <div>9% .</div> </div>
2	C	4	<div> <div>75%</div> <div>25%</div> </div>
2	D	4	<div> <div>75%</div> <div>25%</div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7137 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 Sugar ABC transporter, periplasmic sugar-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	2	0
			3194	2051	557	576	10			
1	B	414	Total	C	N	O	S	0	2	0
			3176	2040	553	573	10			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q53W80
A	174	ARG	LYS	engineered mutation	UNP Q53W80
A	175	THR	ASN	engineered mutation	UNP Q53W80
A	176	PRO	SER	engineered mutation	UNP Q53W80
A	?	-	TRP	deletion	UNP Q53W80
A	177	ARG	ASP	engineered mutation	UNP Q53W80
A	178	THR	VAL	engineered mutation	UNP Q53W80
A	416	HIS	-	expression tag	UNP Q53W80
A	417	HIS	-	expression tag	UNP Q53W80
A	418	HIS	-	expression tag	UNP Q53W80
A	419	HIS	-	expression tag	UNP Q53W80
A	420	HIS	-	expression tag	UNP Q53W80
A	421	HIS	-	expression tag	UNP Q53W80
B	0	MET	-	initiating methionine	UNP Q53W80
B	174	ARG	LYS	engineered mutation	UNP Q53W80
B	175	THR	ASN	engineered mutation	UNP Q53W80
B	176	PRO	SER	engineered mutation	UNP Q53W80
B	?	-	TRP	deletion	UNP Q53W80
B	177	ARG	ASP	engineered mutation	UNP Q53W80
B	178	THR	VAL	engineered mutation	UNP Q53W80
B	416	HIS	-	expression tag	UNP Q53W80
B	417	HIS	-	expression tag	UNP Q53W80
B	418	HIS	-	expression tag	UNP Q53W80
B	419	HIS	-	expression tag	UNP Q53W80
B	420	HIS	-	expression tag	UNP Q53W80

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Chain	Residue	Modelled	Actual	Comment	Reference
B	421	HIS	-	expression tag	UNP Q53W80

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

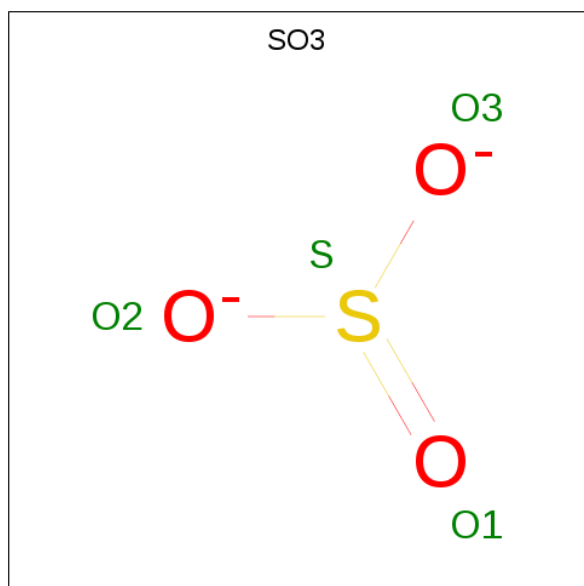


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	4	Total	C	O	0	0	0
			45	24	21			
2	D	4	Total	C	O	0	0	0
			45	24	21			

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	A	2	Total	Cl	0	0
			2	2		

- Molecule 4 is SULFITE ION (three-letter code: SO3) (formula: O₃S).



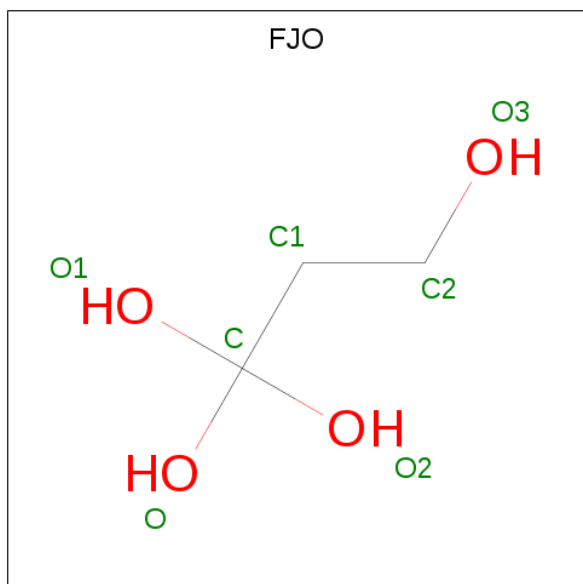
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			4	3	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is propane-1,1,1,3-tetrol (three-letter code: FJO) (formula: $C_3H_8O_4$).

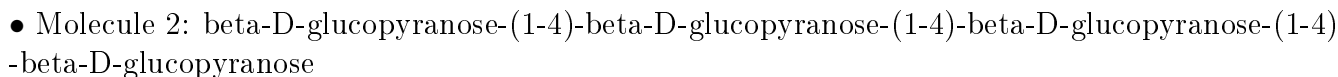
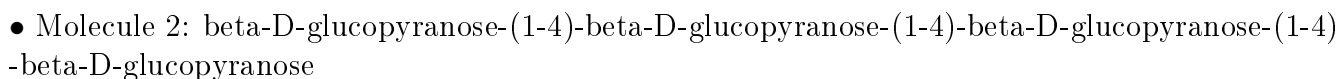


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total 7	C 3	O 4	0	0
6	B	1	Total 7	C 3	O 4	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	336	Total 336	O 336	0	0
7	B	312	Total 312	O 312	0	0

- Molecule 1: Sugar ABC transporter, periplasmic sugar-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.85Å 100.17Å 133.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.53 – 2.05 66.53 – 2.05	Depositor EDS
% Data completeness (in resolution range)	98.3 (66.53-2.05) 98.3 (66.53-2.05)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.24 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.161 , 0.220 0.171 , 0.224	Depositor DCC
R_{free} test set	2467 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.581	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7137	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

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: SO3, BGC, FJO, EDO, CL

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	0.85	2/3289 (0.1%)	0.97	7/4478 (0.2%)
1	B	0.86	4/3269 (0.1%)	0.92	1/4449 (0.0%)
All	All	0.86	6/6558 (0.1%)	0.95	8/8927 (0.1%)

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	161	GLU	CD-OE1	7.67	1.34	1.25
1	A	78	GLU	CD-OE2	-7.34	1.17	1.25
1	B	78	GLU	CD-OE2	-7.30	1.17	1.25
1	B	239	GLU	CD-OE1	6.75	1.33	1.25
1	B	271	GLU	CD-OE2	-5.64	1.19	1.25

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	290	ARG	NE-CZ-NH1	7.28	123.94	120.30
1	B	256	MET	CG-SD-CE	-7.26	88.58	100.20
1	A	326	ARG	NE-CZ-NH1	-6.95	116.82	120.30
1	A	290	ARG	NE-CZ-NH2	-6.71	116.94	120.30
1	A	174	ARG	CG-CD-NE	-6.33	98.50	111.80

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	3194	0	3166	24	0
1	B	3176	0	3157	18	0
2	C	45	0	39	1	0
2	D	45	0	39	3	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
4	A	4	0	0	0	0
5	A	4	0	6	1	0
5	B	4	0	6	0	0
6	A	7	0	0	0	0
6	B	7	0	0	1	0
7	A	336	0	0	6	0
7	B	312	0	0	4	0
All	All	7137	0	6413	42	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.

The worst 5 of 42 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:B:239:GLU:OE1	2:D:1:BGC:O6	1.91	0.88
1:A:401:ASP:HB2	7:A:852:HOH:O	1.77	0.84
1:A:177:ARG:HD2	7:A:719:HOH:O	1.85	0.76
1:B:32:VAL:HG11	1:B:310:LEU:HD21	1.70	0.73
1:A:170:CYS:HB3	1:A:238:ILE:HG23	1.81	0.60

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	416/422 (99%)	413 (99%)	3 (1%)	0	100	100
1	B	414/422 (98%)	412 (100%)	2 (0%)	0	100	100
All	All	830/844 (98%)	825 (99%)	5 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	317/321 (99%)	309 (98%)	8 (2%)	47	40
1	B	315/321 (98%)	304 (96%)	11 (4%)	36	29
All	All	632/642 (98%)	613 (97%)	19 (3%)	41	34

5 of 19 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	9	MET
1	B	33	GLU
1	B	280	VAL
1	A	367	SER
1	B	306	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

8 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	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	BGC	C	1	2	12,12,12	0.86	0	17,17,17	1.62	3 (17%)
2	BGC	C	2	2	11,11,12	0.44	0	15,15,17	1.31	3 (20%)
2	BGC	C	3	2	11,11,12	0.64	0	15,15,17	1.18	2 (13%)
2	BGC	C	4	2	11,11,12	0.94	0	15,15,17	1.32	3 (20%)
2	BGC	D	1	2	12,12,12	0.69	0	17,17,17	0.98	0
2	BGC	D	2	2	11,11,12	0.68	0	15,15,17	1.35	3 (20%)
2	BGC	D	3	2	11,11,12	0.72	0	15,15,17	0.79	0
2	BGC	D	4	2	11,11,12	0.58	0	15,15,17	1.27	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	C	1	2	-	2/2/22/22	0/1/1/1
2	BGC	C	2	2	-	1/2/19/22	0/1/1/1
2	BGC	C	3	2	-	0/2/19/22	0/1/1/1
2	BGC	C	4	2	-	0/2/19/22	0/1/1/1
2	BGC	D	1	2	-	1/2/22/22	0/1/1/1
2	BGC	D	2	2	-	0/2/19/22	0/1/1/1
2	BGC	D	3	2	-	0/2/19/22	0/1/1/1
2	BGC	D	4	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	BGC	C3-C4-C5	5.23	119.56	110.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4	BGC	O5-C5-C6	3.45	112.62	107.20
2	C	3	BGC	O2-C2-C1	3.18	115.65	109.15
2	D	2	BGC	C6-C5-C4	2.64	119.18	113.00
2	C	2	BGC	C6-C5-C4	2.48	118.82	113.00

There are no chirality outliers.

All (4) torsion outliers are listed below:

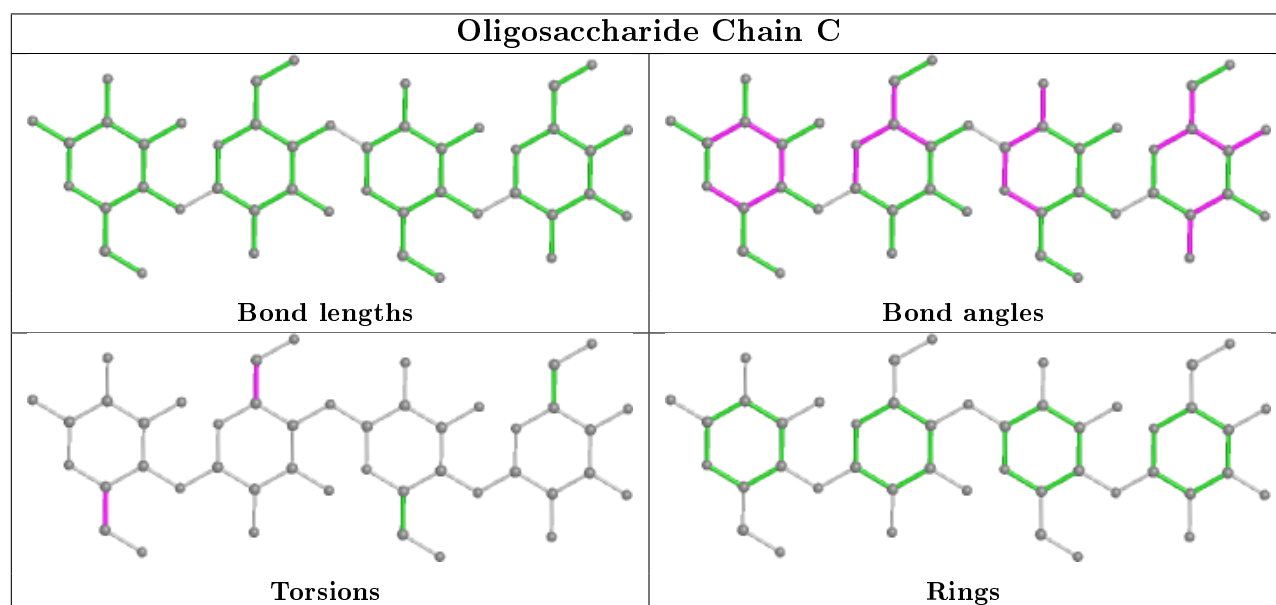
Mol	Chain	Res	Type	Atoms
2	C	1	BGC	C4-C5-C6-O6
2	C	1	BGC	O5-C5-C6-O6
2	D	1	BGC	C4-C5-C6-O6
2	C	2	BGC	C4-C5-C6-O6

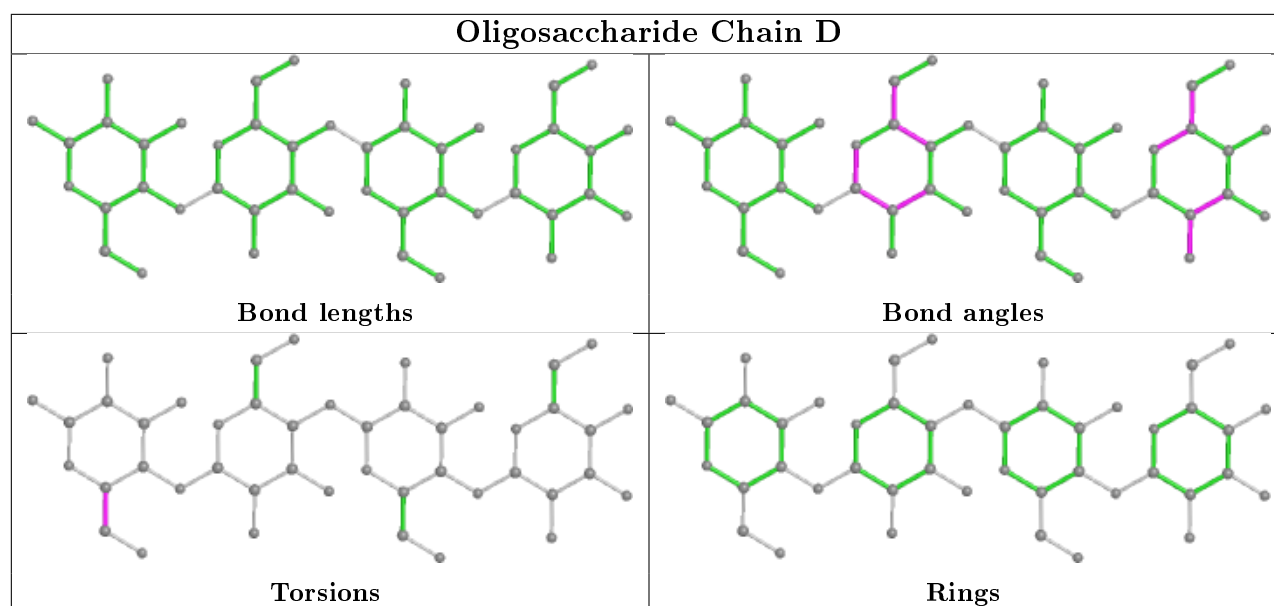
There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	3	BGC	1	0
2	D	2	BGC	1	0
2	C	3	BGC	1	0
2	D	1	BGC	1	0

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 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	$\# Z > 2$
5	EDO	A	504	-	3,3,3	0.23	0	2,2,2	0.21	0
5	EDO	B	502	-	3,3,3	0.37	0	2,2,2	0.46	0
4	SO3	A	503	-	1,3,3	1.18	0	0,3,3	0.00	-
6	FJO	B	503	-	3,6,6	1.07	0	4,8,8	1.12	0
6	FJO	A	505	-	3,6,6	1.62	1 (33%)	4,8,8	1.15	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
5	EDO	A	504	-	-	1/1/1/1	-
5	EDO	B	502	-	-	0/1/1/1	-
6	FJO	B	503	-	-	1/1/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	FJO	A	505	-	-	0/1/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	505	FJO	C1-C2	2.04	1.56	1.52

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	503	FJO	C-C1-C2-O3
5	A	504	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	504	EDO	1	0
6	B	503	FJO	1	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, 95th 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	A	416/422 (98%)	-0.57	1 (0%) 95 95	14, 23, 43, 72	0
1	B	414/422 (98%)	-0.58	1 (0%) 95 95	13, 23, 45, 77	0
All	All	830/844 (98%)	-0.57	2 (0%) 95 95	13, 23, 44, 77	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	393	ARG	2.2
1	A	175	THR	2.2

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](#)

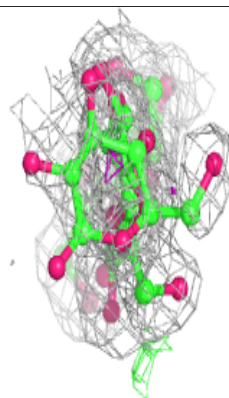
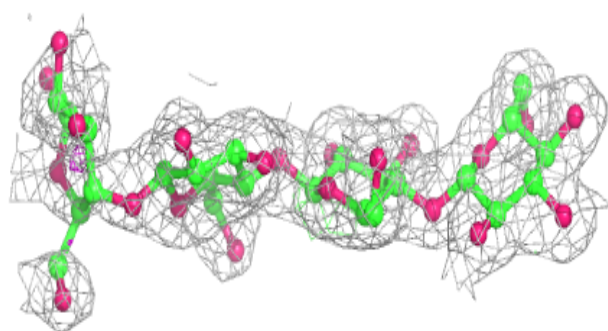
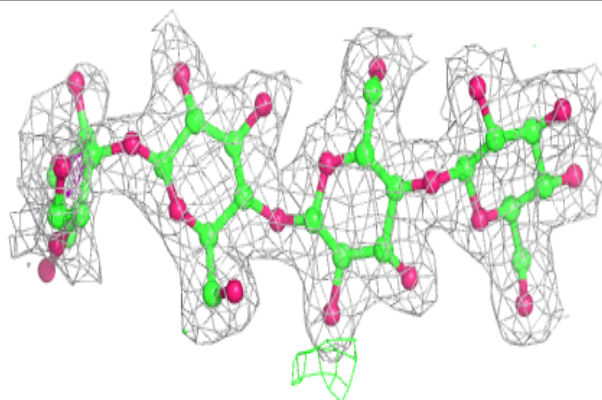
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BGC	C	1	12/12	0.78	0.28	41,55,66,71	0
2	BGC	D	2	11/12	0.95	0.10	15,18,23,33	0
2	BGC	D	1	12/12	0.95	0.13	21,27,36,36	0
2	BGC	C	2	11/12	0.95	0.08	21,25,33,39	0
2	BGC	C	4	11/12	0.97	0.08	13,15,17,17	0
2	BGC	C	3	11/12	0.97	0.08	14,17,19,19	0
2	BGC	D	4	11/12	0.98	0.06	13,14,16,17	0
2	BGC	D	3	11/12	0.98	0.07	14,15,16,16	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.

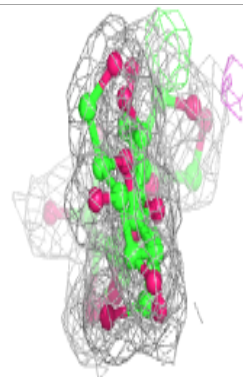
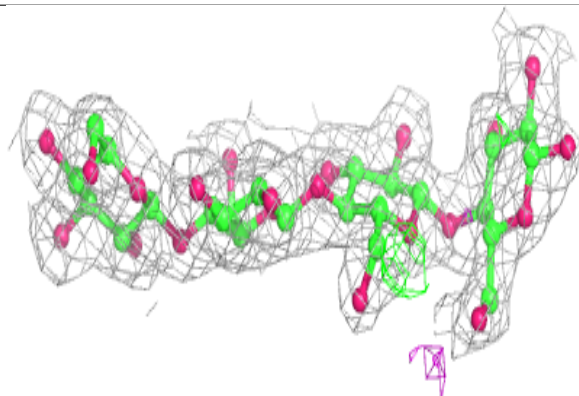
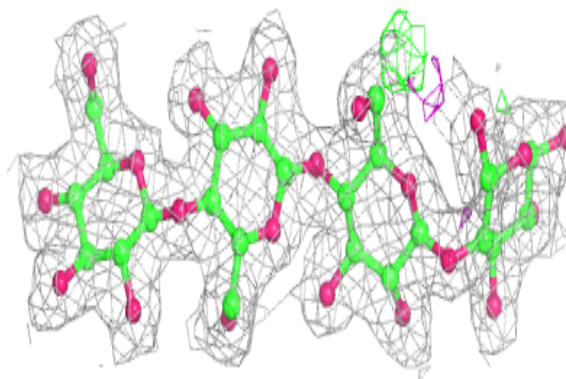
Electron density around Chain C:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain D:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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, 95th 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(\AA^2)	Q<0.9
5	EDO	A	504	4/4	0.83	0.18	38,39,40,43	0
6	FJO	A	505	7/7	0.87	0.16	18,22,28,30	0
5	EDO	B	502	4/4	0.93	0.15	40,41,45,47	0
6	FJO	B	503	7/7	0.94	0.18	23,27,30,31	0
3	CL	A	501	1/1	0.98	0.05	29,29,29,29	0
4	SO3	A	503	4/4	0.98	0.12	41,47,51,53	0
3	CL	B	501	1/1	0.99	0.06	27,27,27,27	0
3	CL	A	502	1/1	0.99	0.07	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.