



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 15, 2020 – 10:42 AM BST

PDB ID : 7C6L
Title : Crystal structure of beta-glycosides-binding protein (W177X) of ABC transporter in a closed state bound to celotriose (Form II)
Authors : Kanaujia, S.P.; Chandravanshi, M.; Samanta, R.
Deposited on : 2020-05-21
Resolution : 2.40 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 ⓘ 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

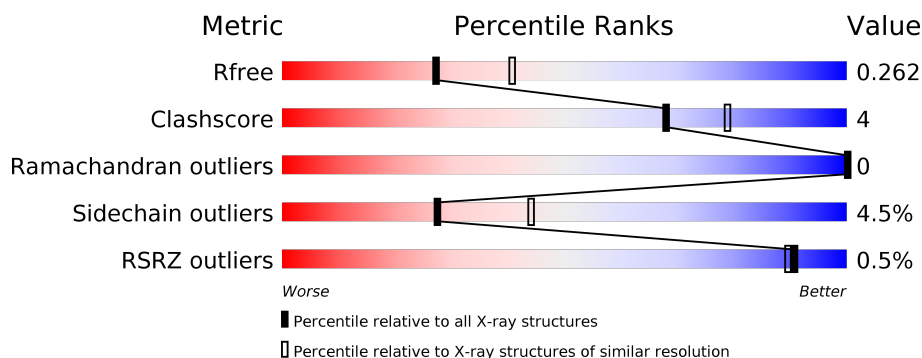
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 style="width: 89%;"></div> <div style="width: 8%;"></div> <div style="width: 3%;"></div> <div style="width: 2%;"></div> </div> <div>89% 8% ..</div>
1	B	422	<div> <div style="width: 86%;"></div> <div style="width: 12%;"></div> <div style="width: 2%;"></div> </div> <div>86% 12% ..</div>
2	C	3	<div> <div style="width: 33%;"></div> <div style="width: 67%;"></div> </div> <div>33% 67%</div>
2	D	3	<div> <div style="width: 67%;"></div> <div style="width: 33%;"></div> </div> <div>67% 33%</div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	EDO	A	505	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6774 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	415	Total	C	N	O	S	0	0	0
			3175	2039	553	573	10			
1	B	416	Total	C	N	O	S	0	0	0
			3185	2045	556	574	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 alpha-L-altropyranose-(1-4)-alpha-D-altropyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	3	Total	C	O	0	0	0
			34	18	16			
2	D	3	Total	C	O	0	0	0
			34	18	16			

- 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	3	Total	Cl	0	0
			3	3		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

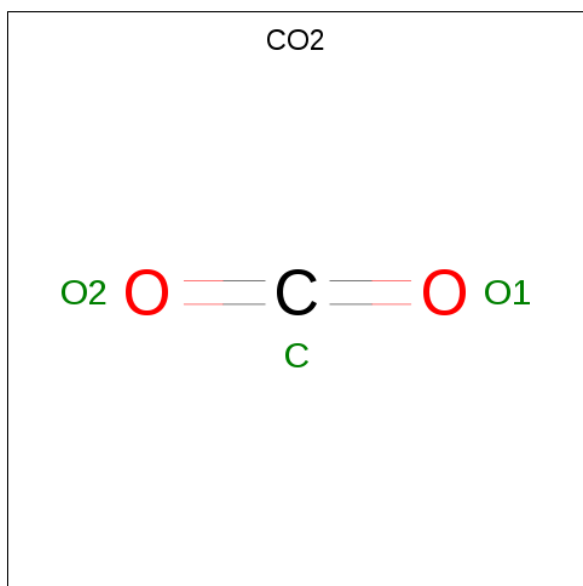
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 6 is CARBON DIOXIDE (three-letter code: CO₂) (formula: CO₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

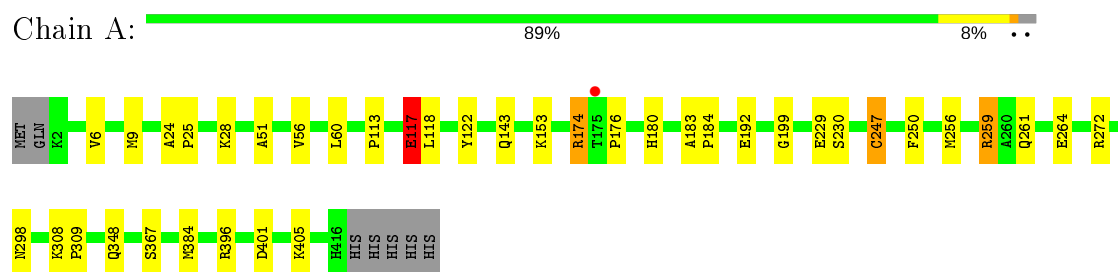
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	178	Total	O	0	0
			178	178		
8	B	149	Total	O	0	0
			149	149		

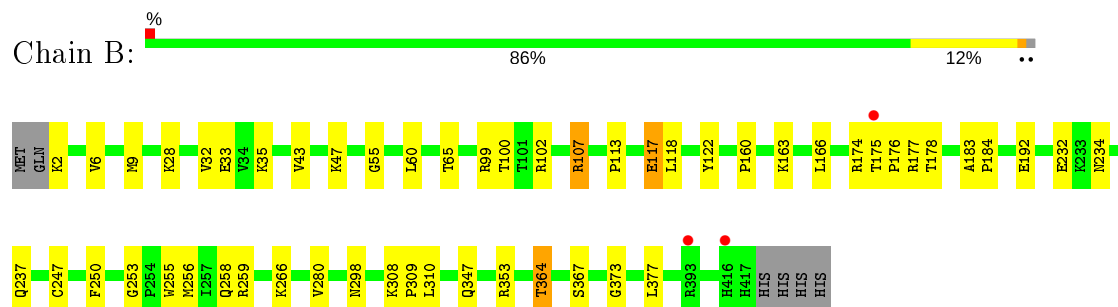
3 Residue-property plots

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: Sugar ABC transporter, periplasmic sugar-binding protein



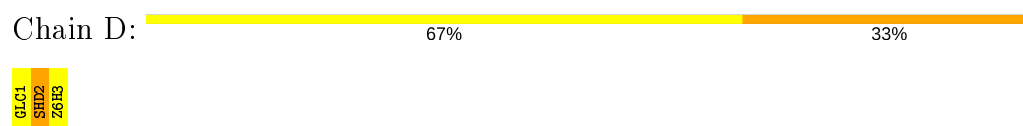
- Molecule 1: Sugar ABC transporter, periplasmic sugar-binding protein



- Molecule 2: alpha-L-altropyranose-(1-4)-alpha-D-altropyranose-(1-4)-alpha-D-glucopyranoside



- Molecule 2: alpha-L-altropyranose-(1-4)-alpha-D-altropyranose-(1-4)-alpha-D-glucopyranoside



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.00Å 100.34Å 133.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.61 – 2.40 66.61 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.8 (66.61-2.40) 99.8 (66.61-2.40)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 2.40Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.191 , 0.263 0.196 , 0.262	Depositor DCC
R_{free} test set	1645 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.479	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6774	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% 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: Z6H, CL, NA, GLC, EDO, CO2, SHD, PEG

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	0.81	0/3263	0.98	6/4442 (0.1%)
1	B	0.78	0/3274	0.92	1/4457 (0.0%)
All	All	0.79	0/6537	0.95	7/8899 (0.1%)

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	A	0	1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	259	ARG	NE-CZ-NH1	6.92	123.76	120.30
1	A	174	ARG	N-CA-CB	6.64	122.55	110.60
1	B	353	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	A	384	MET	CG-SD-CE	5.54	109.06	100.20
1	A	117	GLU	N-CA-CB	-5.52	100.66	110.60
1	A	247	CYS	CB-CA-C	-5.45	99.49	110.40
1	A	230	SER	CB-CA-C	5.16	119.89	110.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	199	GLY	Peptide

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	3175	0	3146	21	0
1	B	3185	0	3153	28	0
2	C	34	0	15	2	0
2	D	34	0	15	1	0
3	A	3	0	0	1	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
5	A	4	0	6	5	0
6	B	3	0	0	0	0
7	B	7	0	10	2	0
8	A	178	0	0	2	0
8	B	149	0	0	3	0
All	All	6774	0	6345	49	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 4.

All (49) 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:174:ARG:O	1:A:176:PRO:HD3	1.74	0.86
1:A:401:ASP:HB2	8:A:733:HOH:O	1.78	0.83
1:B:178:THR:OG1	7:B:503:PEG:H21	1.83	0.79
1:A:396:ARG:HH12	5:A:505:EDO:H12	1.53	0.73
1:A:396:ARG:HH12	5:A:505:EDO:C1	2.04	0.69
1:B:174:ARG:CD	1:B:232:GLU:O	2.42	0.68
1:B:174:ARG:HD3	1:B:232:GLU:O	1.94	0.67
1:B:160:PRO:O	1:B:163:LYS:HE3	1.95	0.66
1:A:122:TYR:CZ	1:A:256:MET:HB2	2.36	0.60
1:B:373:GLY:O	1:B:377:LEU:HG	2.03	0.58
1:A:192:GLU:HG2	3:A:501:CL:CL	2.40	0.58
1:B:107:ARG:CD	1:B:107:ARG:H	2.18	0.57
1:B:99:ARG:HE	1:B:364:THR:HG21	1.70	0.56
1:B:113:PRO:HA	1:B:298:ASN:OD1	2.05	0.56
1:B:160:PRO:O	1:B:163:LYS:CE	2.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:ARG:NE	8:B:607:HOH:O	2.40	0.55
1:B:174:ARG:HD2	1:B:232:GLU:O	2.06	0.54
1:A:113:PRO:HA	1:A:298:ASN:OD1	2.08	0.53
1:B:122:TYR:CZ	1:B:256:MET:HB2	2.43	0.53
1:B:177:ARG:HB2	7:B:503:PEG:H31	1.92	0.52
1:B:183:ALA:N	1:B:184:PRO:CD	2.73	0.51
1:B:234:ASN:OD1	1:B:237:GLN:HG3	2.11	0.50
1:B:55:GLY:HA2	8:B:656:HOH:O	2.11	0.50
1:B:308:LYS:HB2	1:B:309:PRO:HD3	1.94	0.50
1:B:32:VAL:HG11	1:B:310:LEU:HD21	1.94	0.49
1:A:272:ARG:HA	1:A:272:ARG:NE	2.28	0.49
1:A:401:ASP:OD2	1:A:405:LYS:NZ	2.45	0.49
1:B:258:GLN:OE1	1:B:259:ARG:NH1	2.47	0.47
1:A:259:ARG:NH2	2:C:1:GLC:O2	2.47	0.46
1:B:65:THR:HB	1:B:100:THR:HG21	1.98	0.46
1:B:166:LEU:N	1:B:166:LEU:HD12	2.32	0.45
1:B:175:THR:N	1:B:176:PRO:CD	2.81	0.43
1:B:253:GLY:HA3	1:B:255:TRP:CZ3	2.54	0.43
1:B:6:VAL:HA	1:B:60:LEU:O	2.19	0.43
1:A:183:ALA:N	1:A:184:PRO:CD	2.82	0.43
1:A:6:VAL:HA	1:A:60:LEU:O	2.18	0.43
1:A:117:GLU:HG3	2:C:2:SHD:H26	2.01	0.42
1:A:180:HIS:HE1	8:A:625:HOH:O	2.02	0.42
1:B:183:ALA:HB3	1:B:184:PRO:HD3	2.01	0.42
1:A:396:ARG:HH22	5:A:505:EDO:H12	1.85	0.42
1:A:396:ARG:HH12	5:A:505:EDO:H11	1.81	0.42
1:A:51:ALA:HA	1:A:56:VAL:O	2.20	0.42
1:B:117:GLU:HG3	2:D:2:SHD:H26	2.01	0.42
1:A:24:ALA:HB3	1:A:25:PRO:HD3	2.01	0.41
1:A:396:ARG:NH1	5:A:505:EDO:H12	2.28	0.41
1:B:33:GLU:OE2	1:B:35:LYS:HG3	2.20	0.41
1:A:308:LYS:HB2	1:A:309:PRO:HD3	2.03	0.41
1:B:347:GLN:HG3	8:B:708:HOH:O	2.20	0.41
1:A:183:ALA:HB3	1:A:184:PRO:HD3	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	413/422 (98%)	404 (98%)	9 (2%)	0	100	100
1	B	414/422 (98%)	404 (98%)	10 (2%)	0	100	100
All	All	827/844 (98%)	808 (98%)	19 (2%)	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	314/321 (98%)	301 (96%)	13 (4%)	30	48
1	B	315/321 (98%)	300 (95%)	15 (5%)	25	41
All	All	629/642 (98%)	601 (96%)	28 (4%)	27	44

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	MET
1	A	28	LYS
1	A	117	GLU
1	A	118	LEU
1	A	143	GLN
1	A	153	LYS
1	A	229	GLU
1	A	247	CYS

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Mol	Chain	Res	Type
1	A	250	PHE
1	A	261	GLN
1	A	264	GLU
1	A	348	GLN
1	A	367	SER
1	B	2	LYS
1	B	9	MET
1	B	28	LYS
1	B	43	VAL
1	B	47	LYS
1	B	107	ARG
1	B	117	GLU
1	B	118	LEU
1	B	192	GLU
1	B	247	CYS
1	B	250	PHE
1	B	266	LYS
1	B	280	VAL
1	B	364	THR
1	B	367	SER

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

Mol	Chain	Res	Type
1	A	180	HIS
1	A	243	GLN

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 ⓘ

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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GLC	C	1	2	12,12,12	0.46	0	17,17,17	1.40	2 (11%)
2	SHD	C	2	2	11,11,12	0.75	0	15,15,17	1.95	5 (33%)
2	Z6H	C	3	2	11,11,12	0.87	1 (9%)	15,15,17	1.77	3 (20%)
2	GLC	D	1	2	12,12,12	0.60	0	17,17,17	1.33	1 (5%)
2	SHD	D	2	2	11,11,12	0.67	0	15,15,17	1.65	3 (20%)
2	Z6H	D	3	2	11,11,12	0.46	0	15,15,17	1.32	1 (6%)

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	C	1	2	-	0/2/22/22	0/1/1/1
2	SHD	C	2	2	-	1/2/19/22	0/1/1/1
2	Z6H	C	3	2	-	2/2/19/22	0/1/1/1
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	SHD	D	2	2	-	0/2/19/22	0/1/1/1
2	Z6H	D	3	2	-	1/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	Z6H	O5-C1	-2.05	1.40	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	SHD	C1-C2-C3	-4.27	104.42	109.67
2	C	3	Z6H	C1-O5-C5	4.11	117.76	112.19
2	D	3	Z6H	C1-O5-C5	3.87	117.43	112.19
2	D	1	GLC	O5-C1-C2	3.52	116.57	110.28
2	D	2	SHD	C1-C2-C3	-3.18	105.75	109.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	SHD	C3-C4-C5	3.16	115.87	110.24
2	C	3	Z6H	O5-C5-C6	-3.04	102.44	107.20
2	C	1	GLC	O4-C4-C3	2.82	116.86	110.35
2	C	3	Z6H	O2-C2-C3	-2.77	104.59	110.14
2	C	2	SHD	O4-C4-C5	-2.34	103.49	109.30
2	C	2	SHD	C1-O5-C5	2.26	115.25	112.19
2	D	2	SHD	C1-O5-C5	2.22	115.20	112.19
2	C	1	GLC	C3-C4-C5	-2.19	106.33	110.24
2	D	2	SHD	O3-C3-C2	2.17	114.15	109.99
2	C	2	SHD	O5-C5-C4	2.16	116.08	110.83

There are no chirality outliers.

All (4) torsion outliers are listed below:

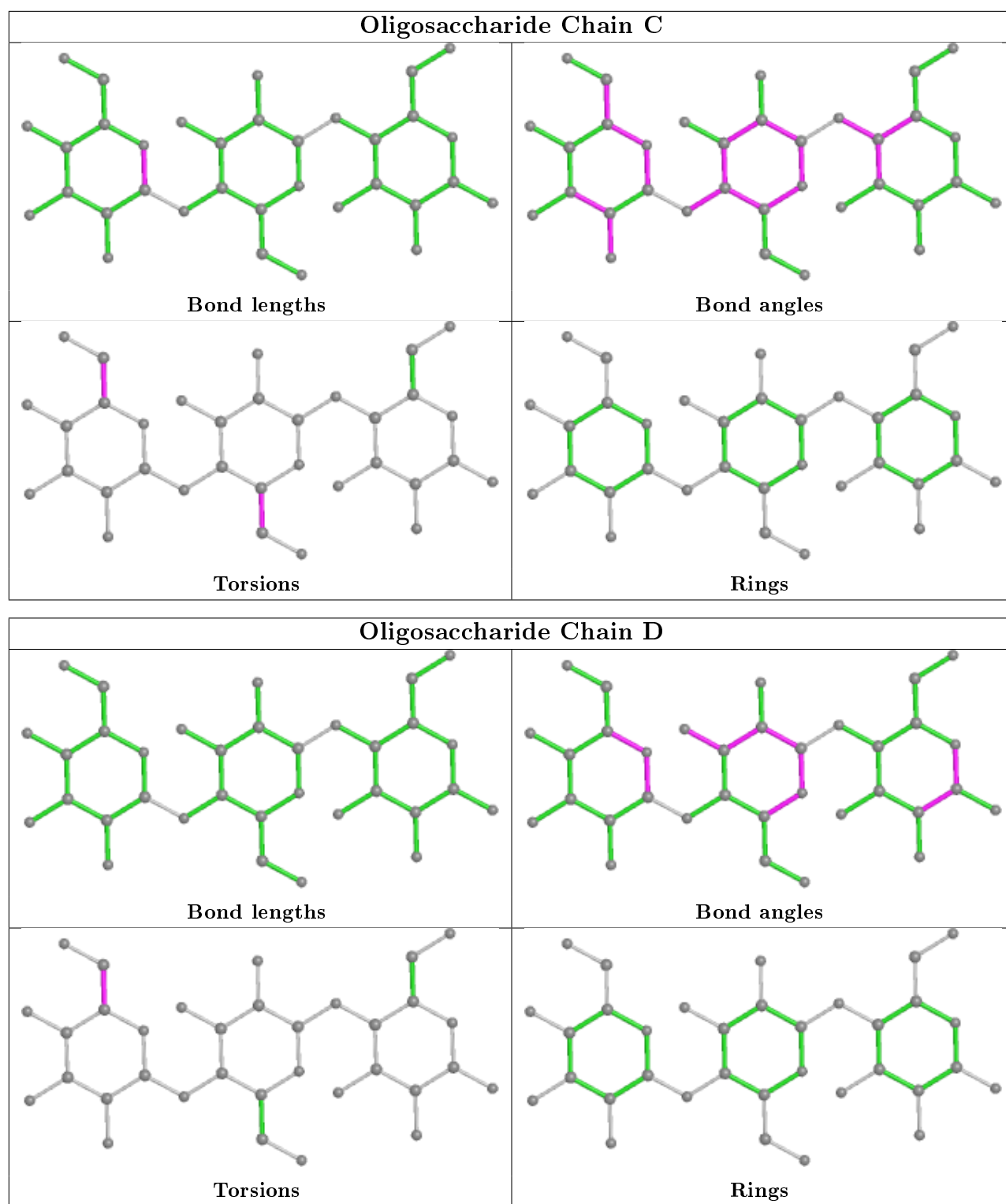
Mol	Chain	Res	Type	Atoms
2	C	3	Z6H	C4-C5-C6-O6
2	C	3	Z6H	O5-C5-C6-O6
2	D	3	Z6H	C4-C5-C6-O6
2	C	2	SHD	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	GLC	1	0
2	D	2	SHD	1	0
2	C	2	SHD	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, 5 are monoatomic - leaving 3 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	505	-	3,3,3	0.30	0	2,2,2	0.54	0
7	PEG	B	503	-	6,6,6	0.27	0	5,5,5	0.22	0
6	CO2	B	502	-	2,2,2	0.05	0	1,1,1	0.63	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	505	-	-	1/1/1/1	-
7	PEG	B	503	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	503	PEG	O1-C1-C2-O2
7	B	503	PEG	O2-C3-C4-O4
5	A	505	EDO	O1-C1-C2-O2
7	B	503	PEG	C1-C2-O2-C3

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	505	EDO	5	0
7	B	503	PEG	2	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	415/422 (98%)	-0.51	1 (0%) 95 94	13, 24, 44, 74	0
1	B	416/422 (98%)	-0.43	3 (0%) 87 86	14, 26, 59, 94	0
All	All	831/844 (98%)	-0.47	4 (0%) 91 89	13, 25, 54, 94	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	175	THR	4.3
1	B	416	HIS	2.8
1	B	393	ARG	2.5
1	B	175	THR	2.0

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](#)

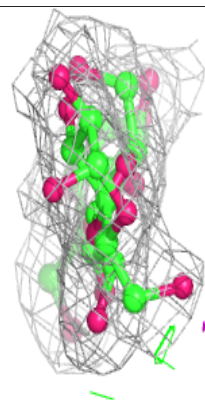
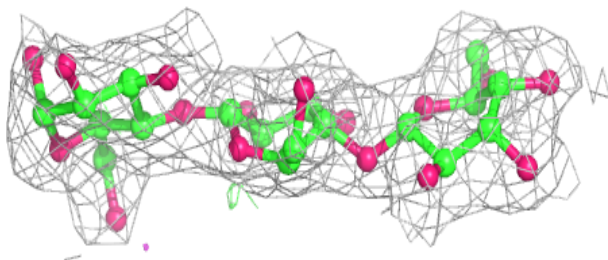
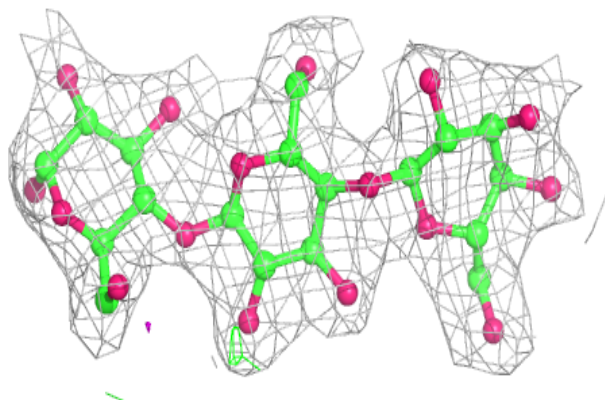
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	GLC	D	1	12/12	0.91	0.12	25,26,28,29	0
2	GLC	C	1	12/12	0.92	0.15	25,31,36,37	0
2	SHD	D	2	11/12	0.95	0.12	23,25,26,26	0
2	Z6H	C	3	11/12	0.95	0.15	18,19,20,23	0
2	Z6H	D	3	11/12	0.95	0.13	18,22,25,25	0
2	SHD	C	2	11/12	0.95	0.10	20,24,26,27	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.

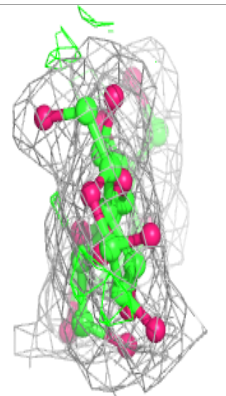
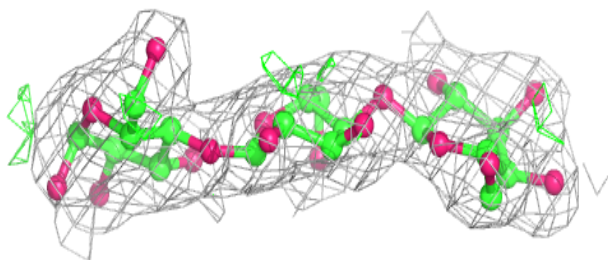
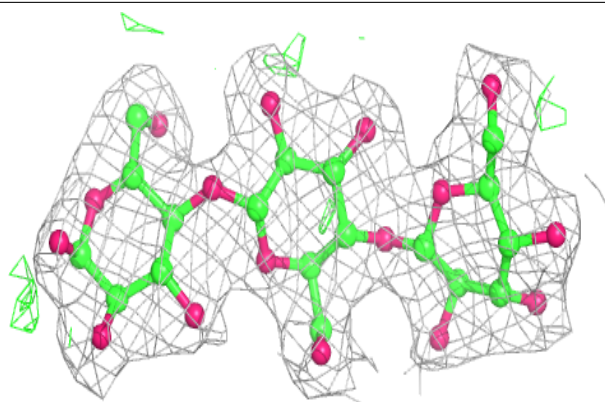
Electron density around Chain C:

$2mF_o-DF_c$ (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	505	4/4	0.76	0.23	31,33,36,38	0
7	PEG	B	503	7/7	0.86	0.16	38,44,46,49	0
6	CO2	B	502	3/3	0.93	0.14	34,34,35,35	0
3	CL	A	501	1/1	0.95	0.09	40,40,40,40	0
4	NA	A	504	1/1	0.96	0.19	30,30,30,30	0
3	CL	A	502	1/1	0.97	0.15	47,47,47,47	0
3	CL	A	503	1/1	0.97	0.08	45,45,45,45	0
3	CL	B	501	1/1	0.98	0.09	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.