



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 10:06 AM UTC

PDB ID : 5D0F / pdb_00005d0f
Title : Crystal Structure of the Candida Glabrata Glycogen Debranching Enzyme (E564Q) in complex with maltopentaose
Authors : Zhai, L.; Xiang, S.
Deposited on : 2015-08-03
Resolution : 3.30 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

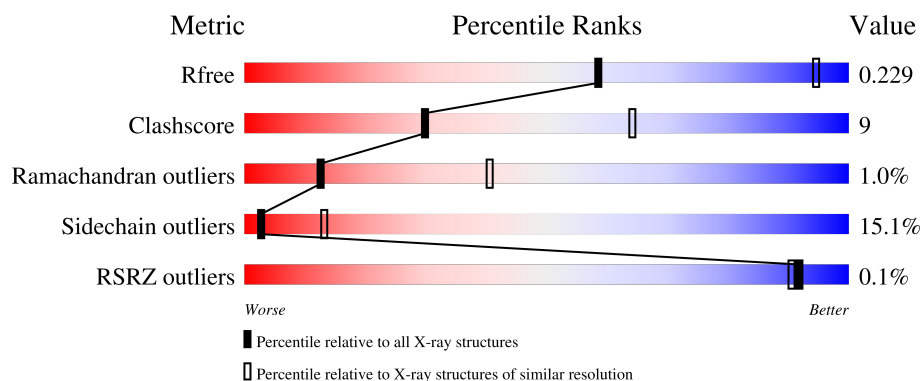
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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 $\leq 5\%$. 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	1528	
1	B	1528	
2	C	4	
2	E	4	
2	H	4	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	I	4	 100%
2	K	4	 100%
3	D	5	 100%
3	G	5	 100%
3	J	5	 100%
3	M	5	 20%60%20%
4	F	3	 100%
4	L	3	 67%33%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 25070 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1526	Total	C	N	O	S	0	0	0
			12278	7830	2066	2330	52			
1	B	1526	Total	C	N	O	S	0	0	0
			12278	7830	2066	2330	52			

There are 2 discrepancies between the modelled and reference sequences:

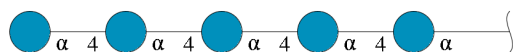
Chain	Residue	Modelled	Actual	Comment	Reference
A	564	GLN	GLU	engineered mutation	UNP Q6FSK0
B	564	GLN	GLU	engineered mutation	UNP Q6FSK0

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



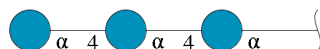
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	4	Total	C	O	0	0	0
			45	24	21			
2	E	4	Total	C	O	0	0	0
			44	24	20			
2	H	4	Total	C	O	0	0	0
			44	24	20			
2	I	4	Total	C	O	0	0	0
			45	24	21			
2	K	4	Total	C	O	0	0	0
			44	24	20			

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	D	5	Total	C	O	0	0	0
			56	30	26			
3	G	5	Total	C	O	0	0	0
			56	30	26			
3	J	5	Total	C	O	0	0	0
			56	30	26			
3	M	5	Total	C	O	0	0	0
			56	30	26			

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

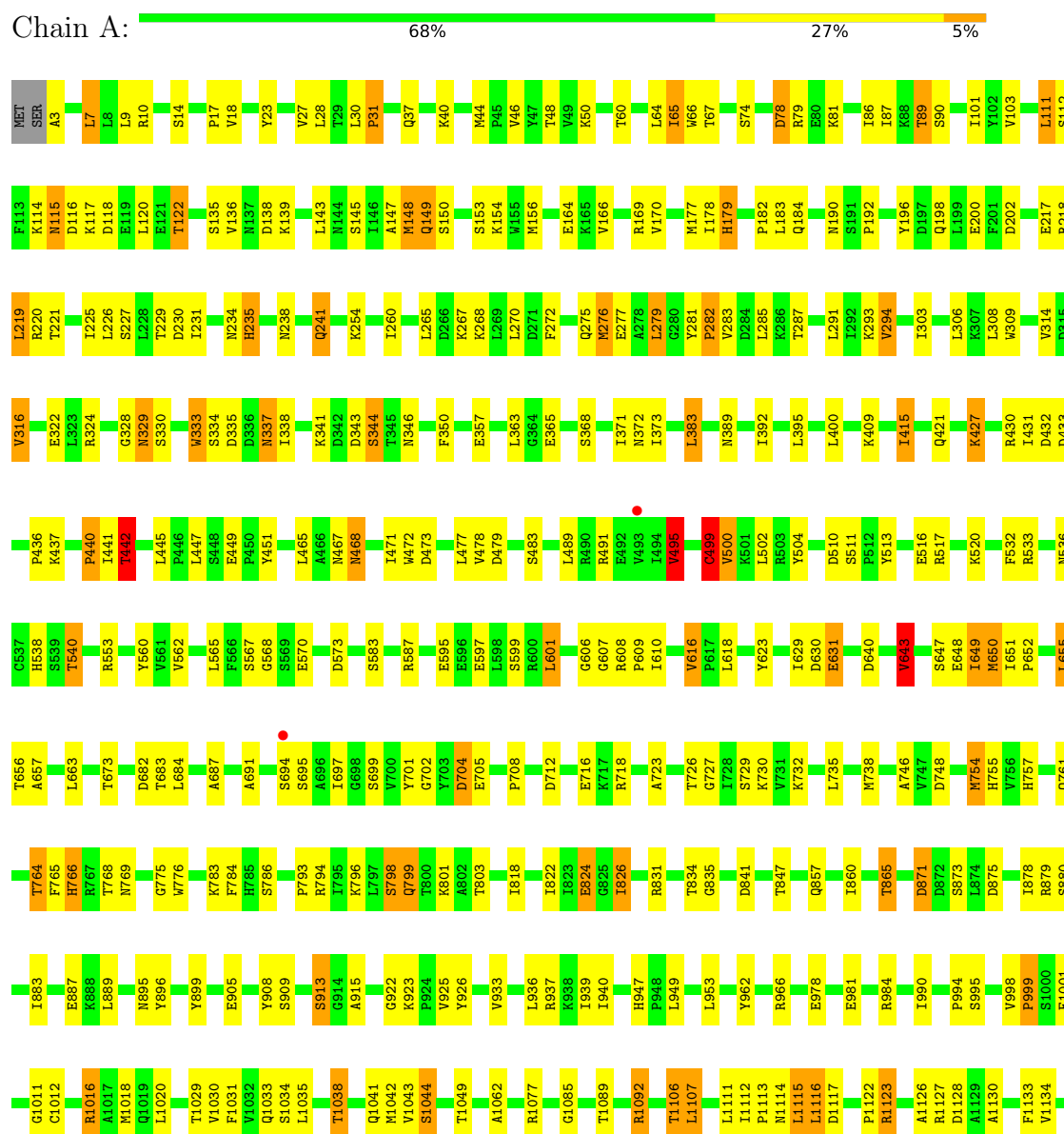


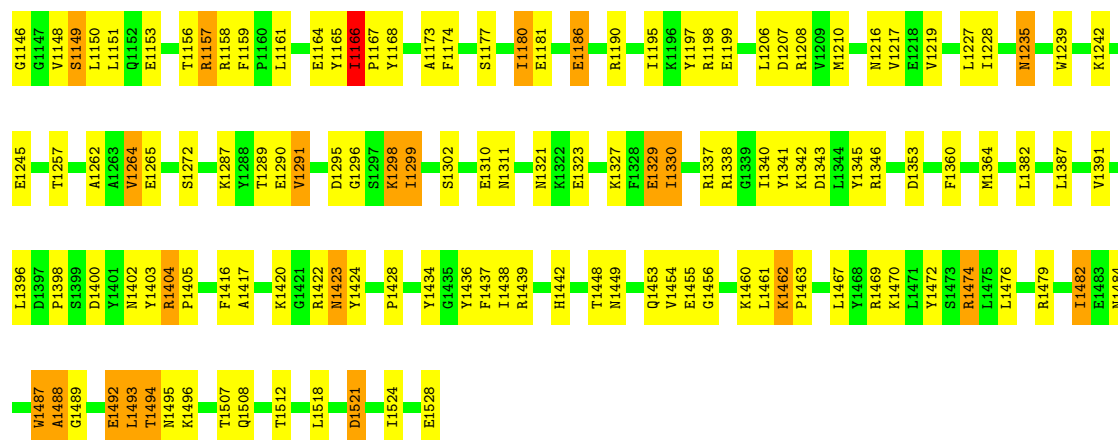
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	F	3	Total	C	O	0	0	0
			34	18	16			
4	L	3	Total	C	O	0	0	0
			34	18	16			

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.

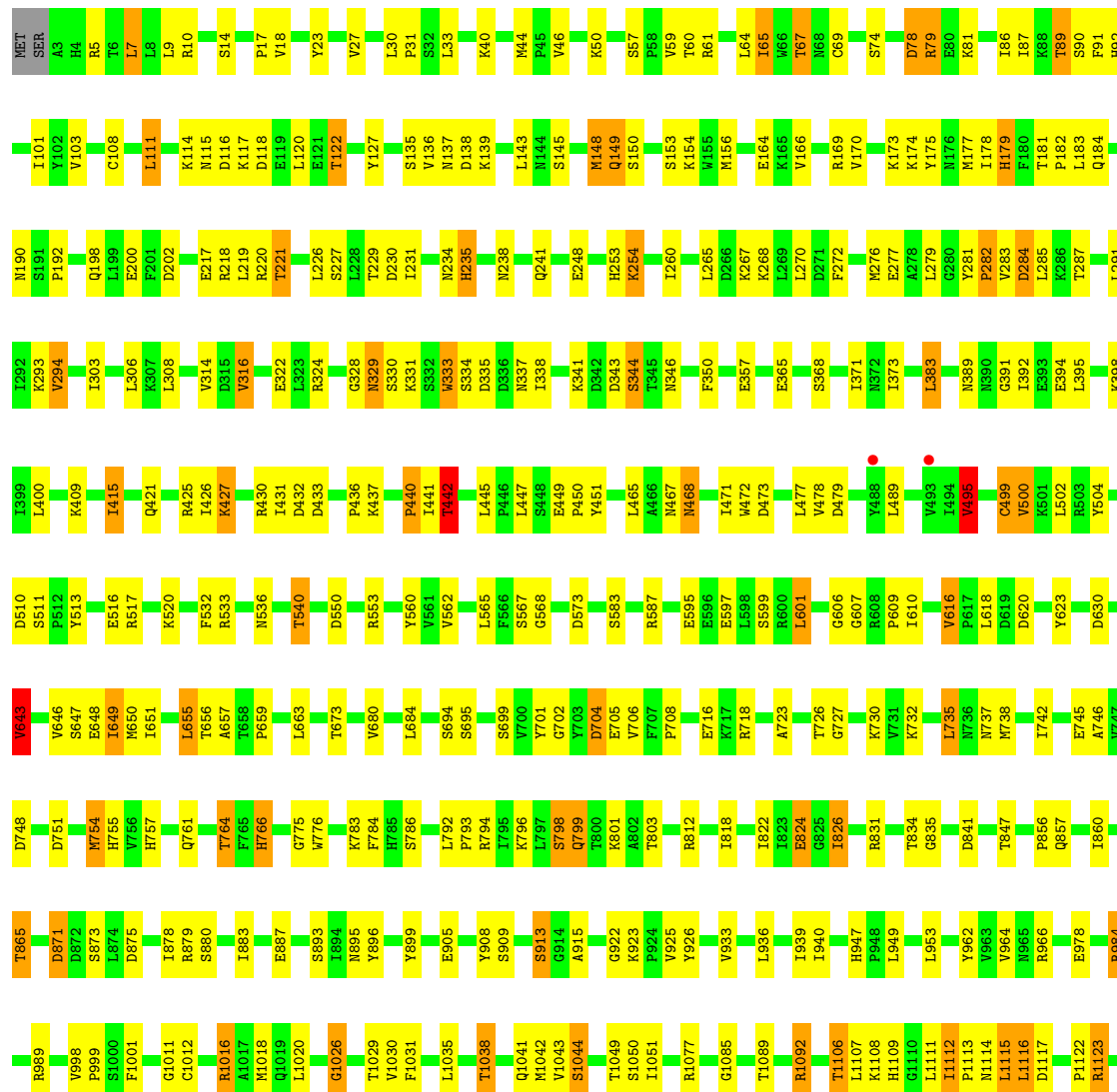
- Molecule 1: Uncharacterized protein

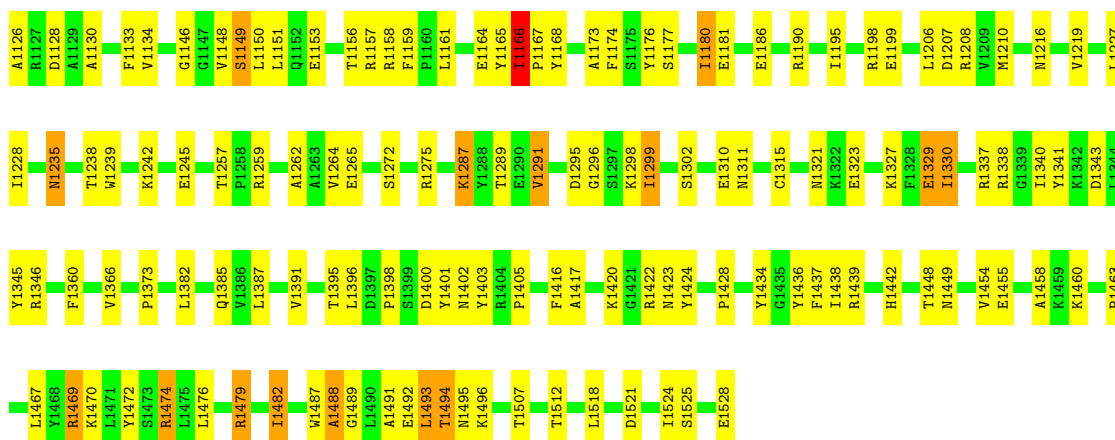




• Molecule 1: Uncharacterized protein

Chain B: 68% 27% 5%





- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain C:  100%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain E:  100%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain H:  100%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain I:  100%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain K:  100%

GLC1
GLC2
GLC3
GLC4

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:  100%

GLC1
GLC2
GLC3
GLC4
GLC5

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain G:  100%

GLC1
GLC2
GLC3
GLC4
GLC5

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain J:  100%

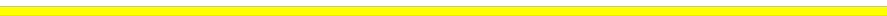
GLC1
GLC2
GLC3
GLC4
GLC5

- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain M:  20% 60% 20%

GLC1
GLC2
GLC3
GLC4
GLC5

- Molecule 4: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain F:  100%

GLC1
GLC2
GLC3

- Molecule 4: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain L:  67% 33%

GLC1
GLC2
GLC3

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.07Å 202.01Å 135.24Å 90.00° 101.32° 90.00°	Depositor
Resolution (Å)	49.05 – 3.30 49.05 – 3.30	Depositor EDS
% Data completeness (in resolution range)	95.8 (49.05-3.30) 95.8 (49.05-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.192 , 0.228 0.197 , 0.229	Depositor DCC
R_{free} test set	3073 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	85.3	Xtriage
Anisotropy	0.512	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 70.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	25070	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.10	9/12589 (0.1%)	1.22	57/17071 (0.3%)
1	B	1.09	14/12589 (0.1%)	1.23	60/17071 (0.4%)
All	All	1.09	23/25178 (0.1%)	1.23	117/34142 (0.3%)

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	3
1	B	0	4
All	All	0	7

The worst 5 of 23 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	812	ARG	CA-C	7.38	1.62	1.52
1	B	122	THR	CA-C	6.69	1.60	1.52
1	B	659	PRO	CA-C	-6.20	1.46	1.52
1	B	655	LEU	C-O	-6.17	1.16	1.23
1	A	695	SER	C-O	-6.12	1.17	1.23

The worst 5 of 117 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	651	ILE	CA-C-N	-10.38	110.24	120.52
1	B	651	ILE	C-N-CA	-10.38	110.24	120.52
1	A	651	ILE	CA-C-N	-8.74	111.87	120.52
1	A	651	ILE	C-N-CA	-8.74	111.87	120.52

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1489	GLY	N-CA-C	8.60	125.98	112.31

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1492	GLU	Peptide
1	A	306	LEU	Peptide
1	A	442	THR	Peptide
1	B	241	GLN	Peptide
1	B	306	LEU	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	12278	0	11962	214	0
1	B	12278	0	11962	208	0
2	C	45	0	39	0	0
2	E	44	0	37	0	0
2	H	44	0	37	0	0
2	I	45	0	39	0	0
2	K	44	0	37	0	0
3	D	56	0	48	0	0
3	G	56	0	48	0	0
3	J	56	0	48	0	0
3	M	56	0	48	1	0
4	F	34	0	30	0	0
4	L	34	0	30	1	0
All	All	25070	0	24365	421	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 9.

The worst 5 of 421 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:1158:ARG:O	1:B:1159:PHE:HD2	1.38	1.04
1:B:1158:ARG:O	1:B:1159:PHE:CD2	2.17	0.97
1:A:1062:ALA:O	1:A:1508:GLN:OE1	1.92	0.86
1:A:230:ASP:OD1	1:A:533:ARG:HD3	1.77	0.84
1:B:230:ASP:OD1	1:B:533:ARG:HD3	1.79	0.83

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	1524/1528 (100%)	1314 (86%)	195 (13%)	15 (1%)	12	40
1	B	1524/1528 (100%)	1321 (87%)	188 (12%)	15 (1%)	12	40
All	All	3048/3056 (100%)	2635 (86%)	383 (13%)	30 (1%)	12	40

5 of 30 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	887	GLU
1	B	887	GLU
1	A	440	PRO
1	A	723	ALA
1	A	1424	TYR

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	1344/1346 (100%)	1139 (85%)	205 (15%)	3	13
1	B	1344/1346 (100%)	1144 (85%)	200 (15%)	3	13
All	All	2688/2692 (100%)	2283 (85%)	405 (15%)	3	13

5 of 405 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	145	SER
1	B	449	GLU
1	B	1476	LEU
1	B	184	GLN
1	B	329	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 44 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	329	ASN
1	B	686	ASN
1	B	346	ASN
1	B	536	ASN
1	B	790	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 ⓘ

46 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	0.86	0	17,17,17	1.38	1 (5%)
2	GLC	C	2	2	11,11,12	0.94	1 (9%)	15,15,17	1.20	1 (6%)
2	GLC	C	3	2	11,11,12	1.20	3 (27%)	15,15,17	1.26	2 (13%)
2	GLC	C	4	2	11,11,12	1.84	3 (27%)	15,15,17	1.37	2 (13%)
3	GLC	D	1	3	12,12,12	0.97	0	17,17,17	1.18	3 (17%)
3	GLC	D	2	3	11,11,12	1.56	2 (18%)	15,15,17	1.43	2 (13%)
3	GLC	D	3	3	11,11,12	0.84	0	15,15,17	1.15	1 (6%)
3	GLC	D	4	3	11,11,12	0.67	0	15,15,17	1.45	3 (20%)
3	GLC	D	5	3	11,11,12	0.77	0	15,15,17	0.85	1 (6%)
2	GLC	E	1	2	11,11,12	1.16	1 (9%)	15,15,17	1.84	5 (33%)
2	GLC	E	2	2	11,11,12	0.87	0	15,15,17	1.07	1 (6%)
2	GLC	E	3	2	11,11,12	0.98	0	15,15,17	1.28	2 (13%)
2	GLC	E	4	2	11,11,12	1.07	1 (9%)	15,15,17	2.08	1 (6%)
4	GLC	F	1	4	12,12,12	0.97	0	17,17,17	0.98	1 (5%)
4	GLC	F	2	4	11,11,12	1.57	2 (18%)	15,15,17	2.65	3 (20%)
4	GLC	F	3	4	11,11,12	0.88	1 (9%)	15,15,17	1.53	1 (6%)
3	GLC	G	1	3	12,12,12	0.96	0	17,17,17	1.00	1 (5%)
3	GLC	G	2	3	11,11,12	1.06	1 (9%)	15,15,17	1.06	1 (6%)
3	GLC	G	3	3	11,11,12	0.86	0	15,15,17	1.14	1 (6%)
3	GLC	G	4	3	11,11,12	0.85	0	15,15,17	1.30	1 (6%)
3	GLC	G	5	3	11,11,12	1.29	2 (18%)	15,15,17	1.59	4 (26%)
2	GLC	H	1	2	11,11,12	1.25	1 (9%)	15,15,17	1.01	1 (6%)
2	GLC	H	2	2	11,11,12	1.57	3 (27%)	15,15,17	1.77	3 (20%)
2	GLC	H	3	2	11,11,12	1.20	2 (18%)	15,15,17	1.88	1 (6%)
2	GLC	H	4	2	11,11,12	0.96	0	15,15,17	1.37	1 (6%)
2	GLC	I	1	2	12,12,12	1.08	0	17,17,17	1.31	2 (11%)
2	GLC	I	2	2	11,11,12	0.90	0	15,15,17	1.75	4 (26%)
2	GLC	I	3	2	11,11,12	1.26	1 (9%)	15,15,17	2.43	2 (13%)
2	GLC	I	4	2	11,11,12	1.62	2 (18%)	15,15,17	1.67	2 (13%)
3	GLC	J	1	3	12,12,12	1.24	2 (16%)	17,17,17	1.04	1 (5%)
3	GLC	J	2	3	11,11,12	2.04	3 (27%)	15,15,17	2.69	1 (6%)
3	GLC	J	3	3	11,11,12	0.81	0	15,15,17	1.47	3 (20%)
3	GLC	J	4	3	11,11,12	0.91	1 (9%)	15,15,17	1.11	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GLC	J	5	3	11,11,12	0.95	1 (9%)	15,15,17	1.23	1 (6%)
2	GLC	K	1	2	11,11,12	1.31	3 (27%)	15,15,17	2.91	6 (40%)
2	GLC	K	2	2	11,11,12	0.82	1 (9%)	15,15,17	1.15	3 (20%)
2	GLC	K	3	2	11,11,12	1.08	1 (9%)	15,15,17	1.62	3 (20%)
2	GLC	K	4	2	11,11,12	1.19	2 (18%)	15,15,17	1.76	1 (6%)
4	GLC	L	1	4	12,12,12	0.78	0	17,17,17	0.87	1 (5%)
4	GLC	L	2	4	11,11,12	0.89	0	15,15,17	1.90	2 (13%)
4	GLC	L	3	4	11,11,12	0.71	0	15,15,17	1.74	1 (6%)
3	GLC	M	1	3	12,12,12	0.72	0	17,17,17	0.88	1 (5%)
3	GLC	M	2	3	11,11,12	0.88	0	15,15,17	0.92	0
3	GLC	M	3	3	11,11,12	0.85	0	15,15,17	0.90	1 (6%)
3	GLC	M	4	3	11,11,12	0.85	0	15,15,17	1.62	1 (6%)
3	GLC	M	5	3	11,11,12	1.03	0	15,15,17	1.21	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	C	1	2	-	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	GLC	C	3	2	-	0/2/19/22	0/1/1/1
2	GLC	C	4	2	-	1/2/19/22	0/1/1/1
3	GLC	D	1	3	-	0/2/22/22	0/1/1/1
3	GLC	D	2	3	-	2/2/19/22	0/1/1/1
3	GLC	D	3	3	-	2/2/19/22	0/1/1/1
3	GLC	D	4	3	-	1/2/19/22	0/1/1/1
3	GLC	D	5	3	-	2/2/19/22	0/1/1/1
2	GLC	E	1	2	-	2/2/19/22	0/1/1/1
2	GLC	E	2	2	-	1/2/19/22	0/1/1/1
2	GLC	E	3	2	-	2/2/19/22	0/1/1/1
2	GLC	E	4	2	-	1/2/19/22	0/1/1/1
4	GLC	F	1	4	-	1/2/22/22	0/1/1/1
4	GLC	F	2	4	-	0/2/19/22	0/1/1/1
4	GLC	F	3	4	-	1/2/19/22	0/1/1/1
3	GLC	G	1	3	-	0/2/22/22	0/1/1/1
3	GLC	G	2	3	-	0/2/19/22	0/1/1/1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GLC	G	3	3	-	0/2/19/22	0/1/1/1
3	GLC	G	4	3	-	0/2/19/22	0/1/1/1
3	GLC	G	5	3	-	0/2/19/22	0/1/1/1
2	GLC	H	1	2	-	2/2/19/22	0/1/1/1
2	GLC	H	2	2	-	2/2/19/22	0/1/1/1
2	GLC	H	3	2	-	0/2/19/22	0/1/1/1
2	GLC	H	4	2	-	2/2/19/22	0/1/1/1
2	GLC	I	1	2	-	2/2/22/22	0/1/1/1
2	GLC	I	2	2	-	0/2/19/22	0/1/1/1
2	GLC	I	3	2	-	2/2/19/22	0/1/1/1
2	GLC	I	4	2	-	1/2/19/22	0/1/1/1
3	GLC	J	1	3	-	2/2/22/22	0/1/1/1
3	GLC	J	2	3	-	1/2/19/22	0/1/1/1
3	GLC	J	3	3	-	2/2/19/22	0/1/1/1
3	GLC	J	4	3	-	2/2/19/22	0/1/1/1
3	GLC	J	5	3	-	2/2/19/22	0/1/1/1
2	GLC	K	1	2	-	0/2/19/22	0/1/1/1
2	GLC	K	2	2	-	2/2/19/22	0/1/1/1
2	GLC	K	3	2	-	0/2/19/22	0/1/1/1
2	GLC	K	4	2	-	1/2/19/22	0/1/1/1
4	GLC	L	1	4	-	0/2/22/22	0/1/1/1
4	GLC	L	2	4	-	1/2/19/22	0/1/1/1
4	GLC	L	3	4	-	2/2/19/22	0/1/1/1
3	GLC	M	1	3	-	2/2/22/22	0/1/1/1
3	GLC	M	2	3	-	0/2/19/22	0/1/1/1
3	GLC	M	3	3	-	2/2/19/22	0/1/1/1
3	GLC	M	4	3	-	1/2/19/22	0/1/1/1
3	GLC	M	5	3	-	0/2/19/22	0/1/1/1

The worst 5 of 40 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	2	GLC	O5-C1	4.05	1.50	1.43
3	J	2	GLC	O5-C5	4.03	1.51	1.43
2	I	4	GLC	O5-C1	3.86	1.50	1.43
4	F	2	GLC	O5-C5	3.71	1.50	1.43
2	C	4	GLC	O5-C5	3.67	1.50	1.43

The worst 5 of 84 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	J	2	GLC	C1-O5-C5	9.71	125.20	112.19
4	F	2	GLC	C1-O5-C5	9.02	124.28	112.19
2	I	3	GLC	C1-O5-C5	8.64	123.77	112.19
2	E	4	GLC	C1-O5-C5	7.08	121.67	112.19
2	K	1	GLC	C1-O5-C5	6.72	121.20	112.19

There are no chirality outliers.

5 of 49 torsion outliers are listed below:

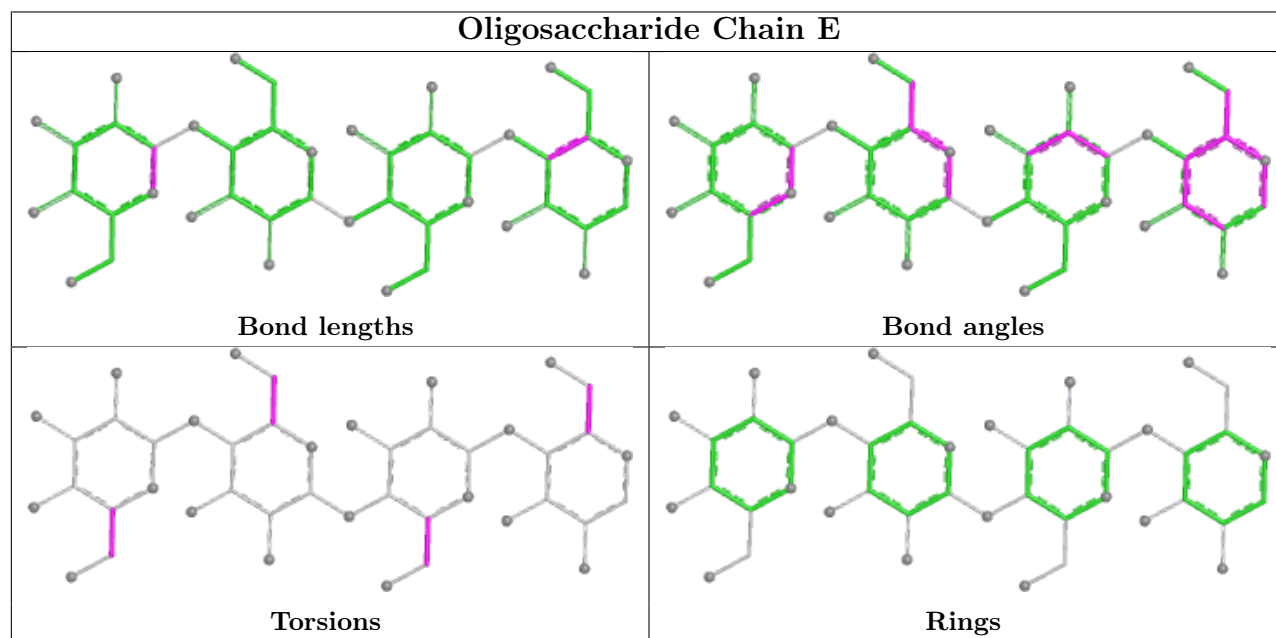
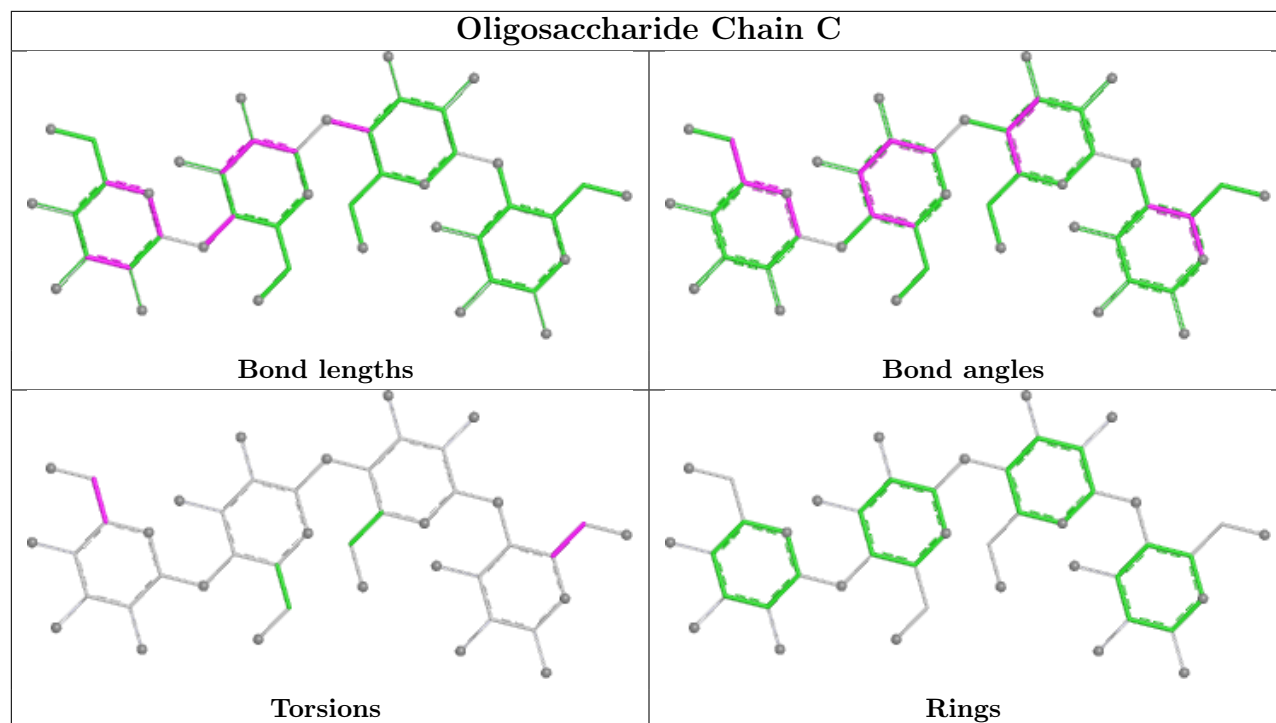
Mol	Chain	Res	Type	Atoms
2	H	2	GLC	C4-C5-C6-O6
2	H	4	GLC	O5-C5-C6-O6
3	J	3	GLC	O5-C5-C6-O6
3	D	3	GLC	O5-C5-C6-O6
2	I	3	GLC	O5-C5-C6-O6

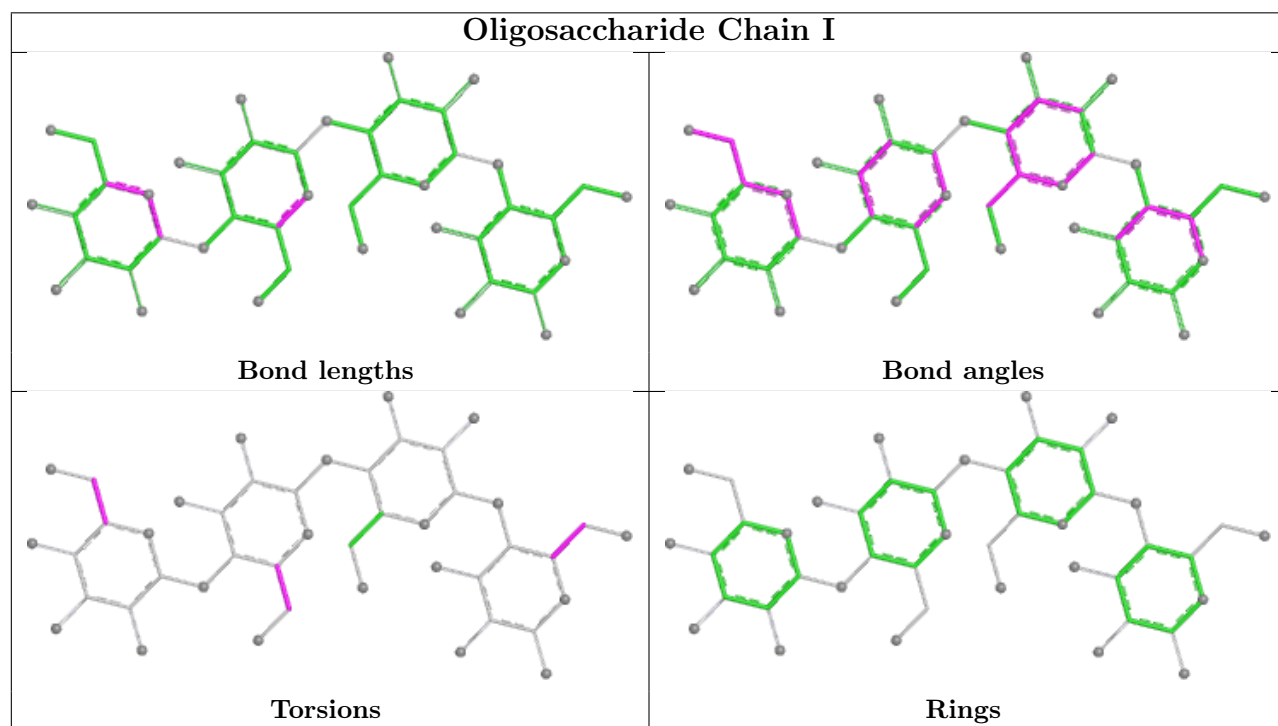
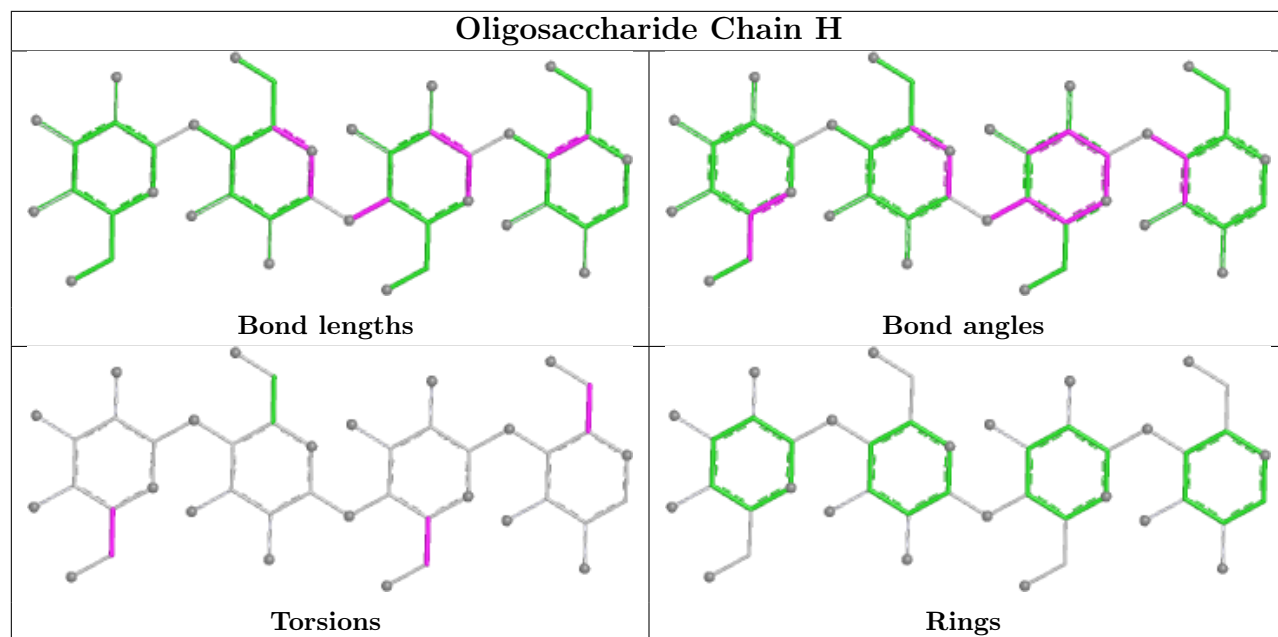
There are no ring outliers.

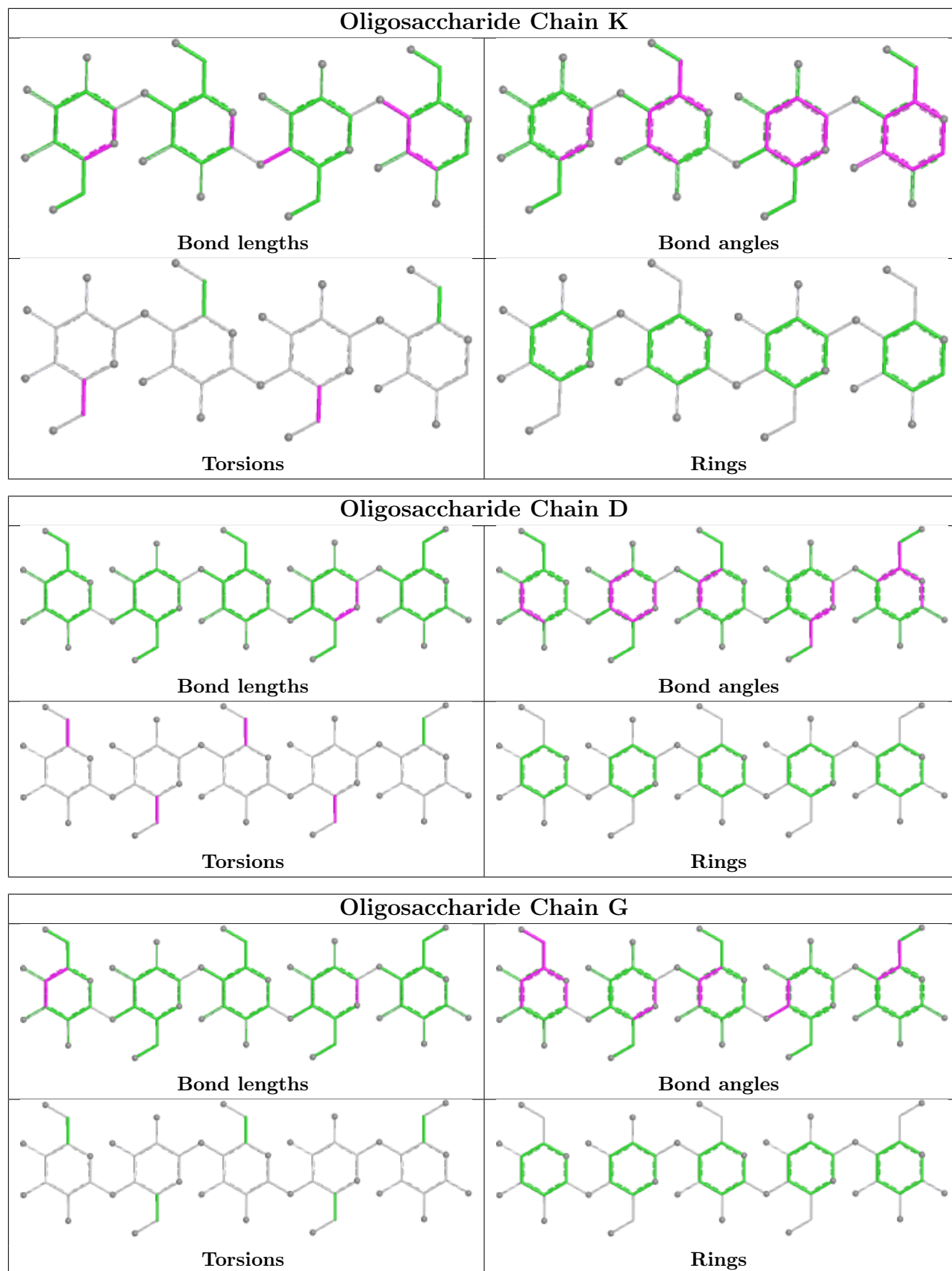
2 monomers are involved in 2 short contacts:

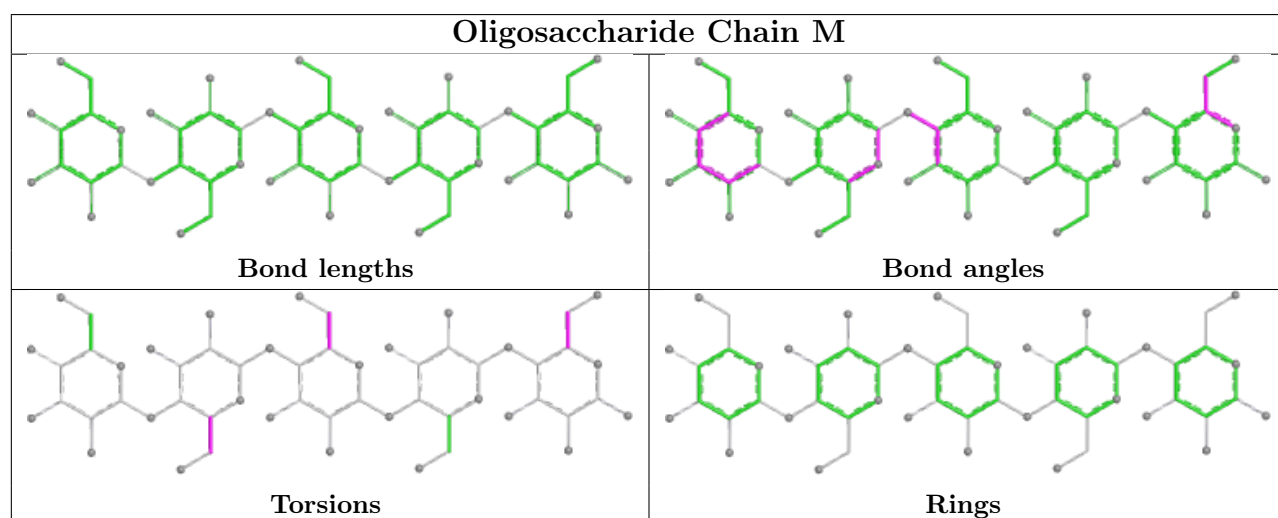
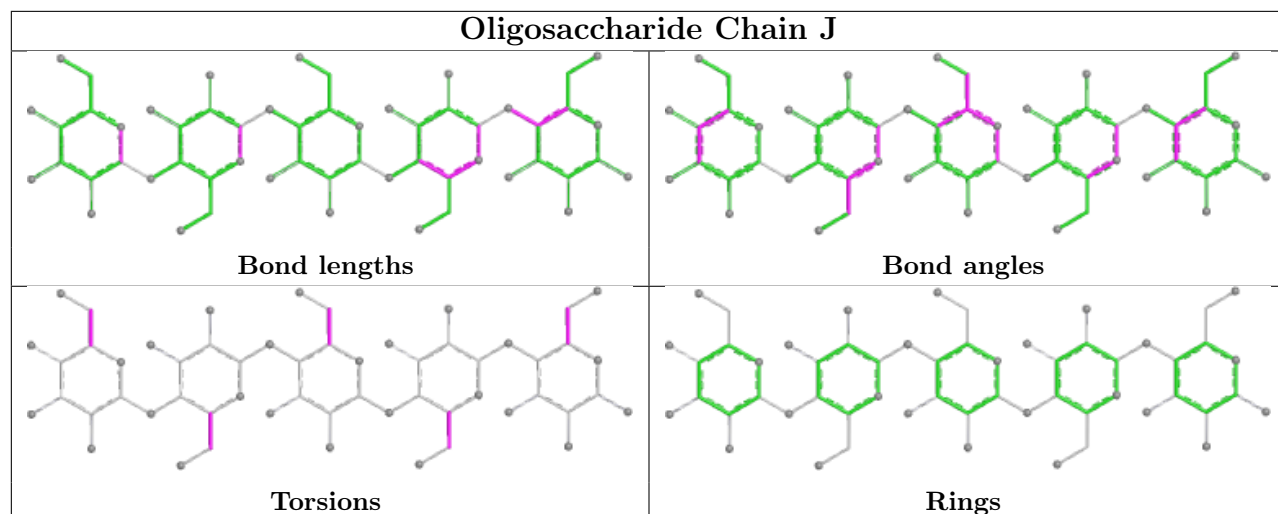
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	1	GLC	1	0
3	M	4	GLC	1	0

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

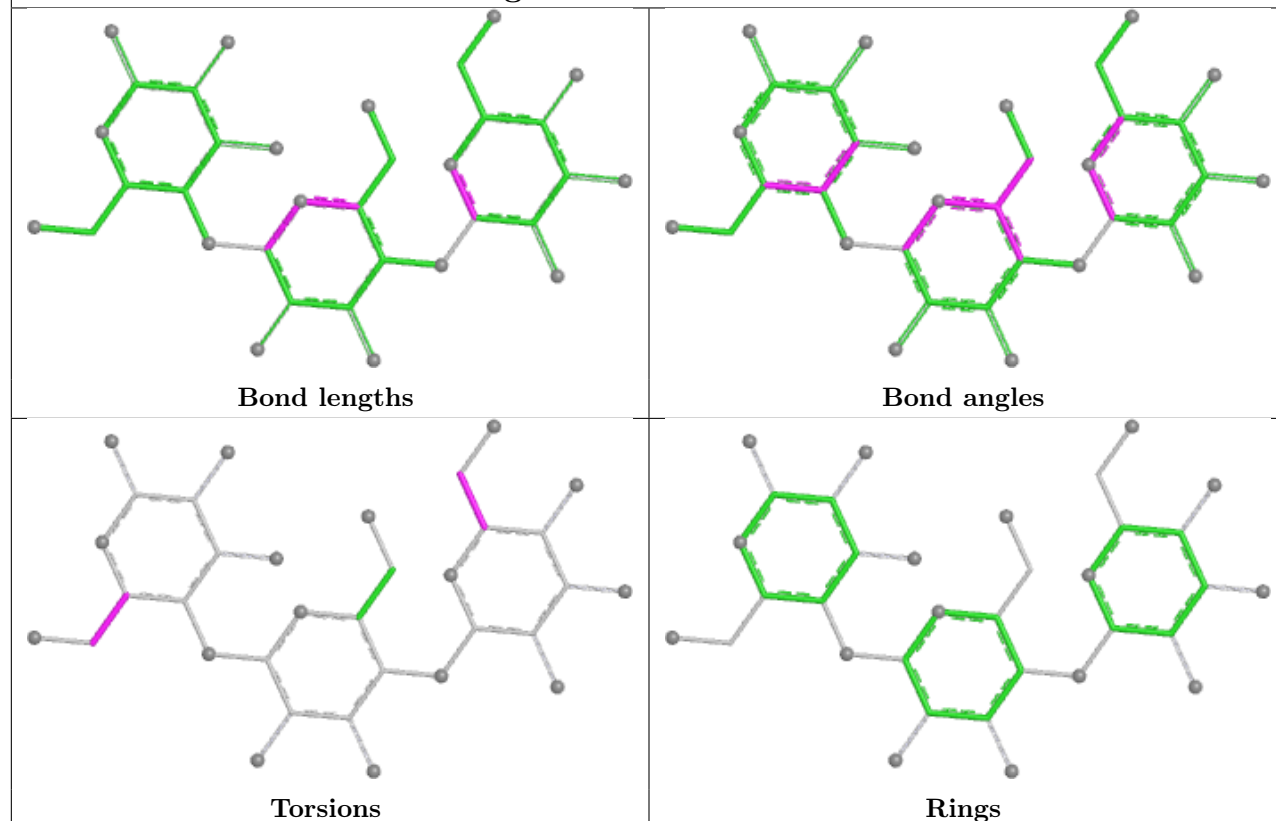




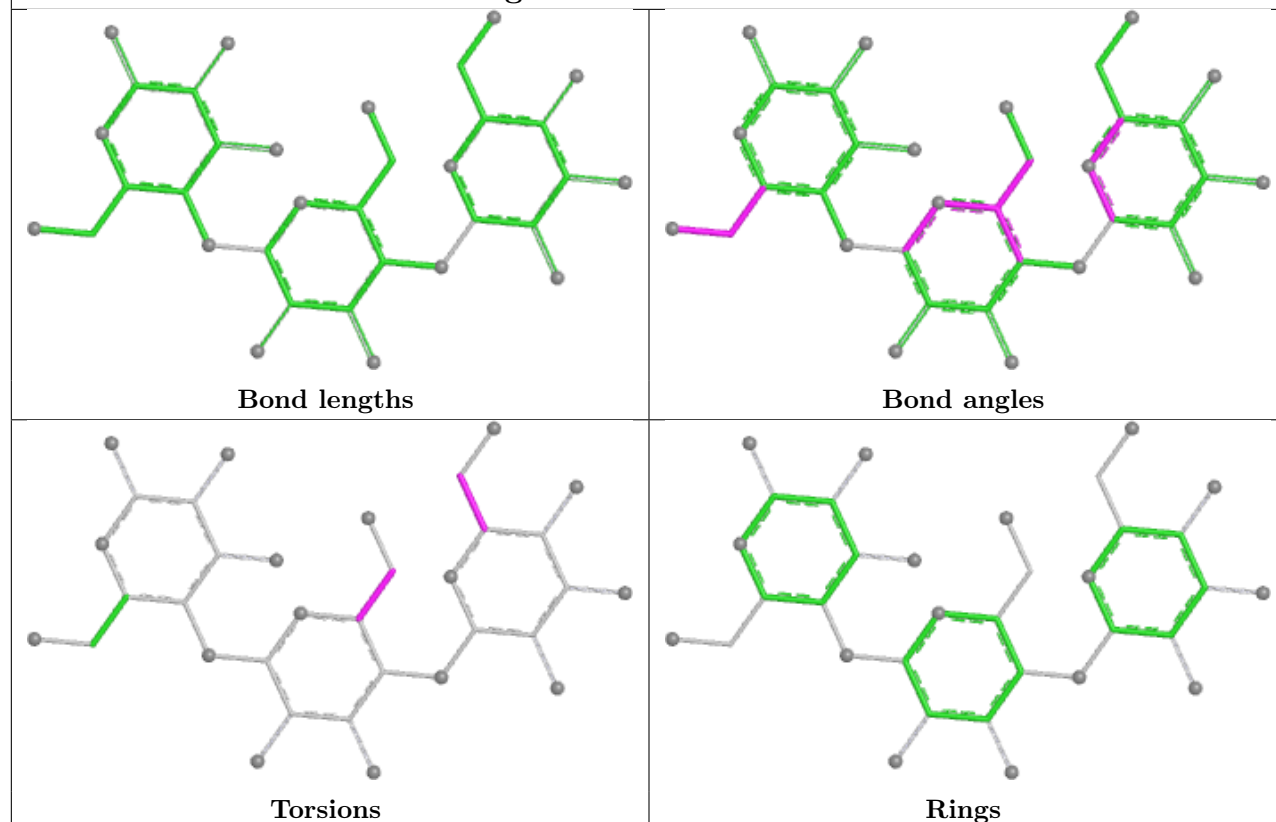




Oligosaccharide Chain F



Oligosaccharide Chain L



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](#)

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	1526/1528 (99%)	-0.45	2 (0%)	92 90	50, 86, 144, 176	0
1	B	1526/1528 (99%)	-0.45	2 (0%)	92 90	50, 83, 146, 190	0
All	All	3052/3056 (99%)	-0.45	4 (0%)	92 90	50, 85, 145, 190	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	493	VAL	2.4
1	A	493	VAL	2.3
1	A	694	SER	2.2
1	B	488	TYR	2.1

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
3	GLC	M	2	11/12	0.49	0.11	150,177,181,188	0
3	GLC	M	5	11/12	0.50	0.12	186,201,208,211	0
2	GLC	H	1	11/12	0.54	0.10	113,137,147,147	0
3	GLC	D	5	11/12	0.54	0.10	144,152,172,173	0
2	GLC	H	4	11/12	0.57	0.07	154,180,189,196	0

Continued on next page...

Continued from previous page...

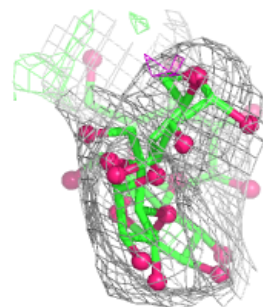
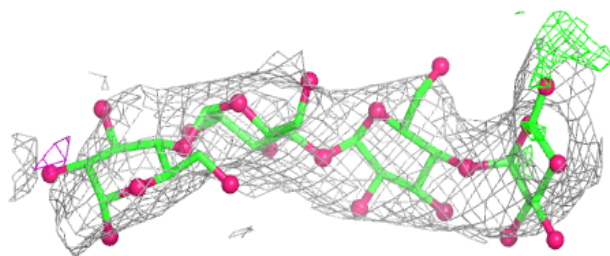
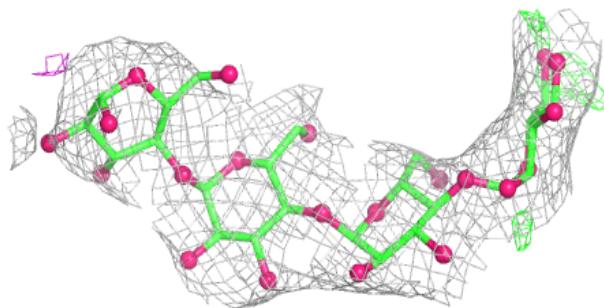
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GLC	F	1	12/12	0.63	0.09	157,161,164,164	0
4	GLC	L	3	11/12	0.65	0.08	130,145,159,162	0
4	GLC	F	3	11/12	0.70	0.07	152,169,173,179	0
3	GLC	J	5	11/12	0.71	0.09	121,129,135,137	0
2	GLC	H	3	11/12	0.71	0.09	147,167,186,188	0
2	GLC	H	2	11/12	0.72	0.11	135,152,165,172	0
3	GLC	G	1	12/12	0.73	0.13	131,183,190,194	0
2	GLC	K	1	11/12	0.76	0.09	104,121,127,130	0
4	GLC	F	2	11/12	0.77	0.08	129,167,181,185	0
3	GLC	J	1	12/12	0.78	0.13	129,138,144,145	0
2	GLC	C	4	11/12	0.78	0.13	98,128,136,141	0
3	GLC	M	1	12/12	0.78	0.10	137,168,174,178	0
3	GLC	M	3	11/12	0.79	0.10	146,173,184,185	0
3	GLC	G	5	11/12	0.79	0.08	174,185,191,192	0
3	GLC	J	2	11/12	0.80	0.13	123,129,144,145	0
2	GLC	I	4	11/12	0.81	0.10	110,127,136,153	0
2	GLC	E	1	11/12	0.81	0.09	120,125,132,135	0
4	GLC	L	2	11/12	0.82	0.08	124,134,166,190	0
3	GLC	D	2	11/12	0.82	0.10	126,132,138,142	0
3	GLC	J	3	11/12	0.84	0.09	105,117,120,125	0
3	GLC	J	4	11/12	0.84	0.09	108,120,128,128	0
2	GLC	K	4	11/12	0.84	0.10	108,120,135,136	0
2	GLC	I	1	12/12	0.85	0.12	138,144,158,162	0
3	GLC	D	1	12/12	0.85	0.12	117,135,142,144	0
2	GLC	C	1	12/12	0.86	0.12	124,143,157,171	0
3	GLC	G	2	11/12	0.87	0.08	138,157,166,171	0
3	GLC	M	4	11/12	0.87	0.09	159,174,189,199	0
4	GLC	L	1	12/12	0.87	0.06	114,131,133,141	0
2	GLC	E	2	11/12	0.87	0.09	92,108,117,118	0
2	GLC	E	4	11/12	0.87	0.08	105,119,126,132	0
3	GLC	G	3	11/12	0.88	0.08	151,166,173,187	0
3	GLC	G	4	11/12	0.91	0.10	147,197,208,208	0
2	GLC	I	3	11/12	0.91	0.10	105,118,134,142	0
3	GLC	D	3	11/12	0.91	0.08	112,129,140,143	0
2	GLC	I	2	11/12	0.92	0.07	102,109,115,120	0
2	GLC	C	3	11/12	0.92	0.08	107,113,133,135	0
3	GLC	D	4	11/12	0.93	0.07	123,133,142,148	0
2	GLC	K	3	11/12	0.94	0.08	91,93,100,105	0
2	GLC	E	3	11/12	0.95	0.07	71,78,90,91	0
2	GLC	K	2	11/12	0.95	0.07	112,113,119,119	0
2	GLC	C	2	11/12	0.96	0.09	102,114,121,128	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.

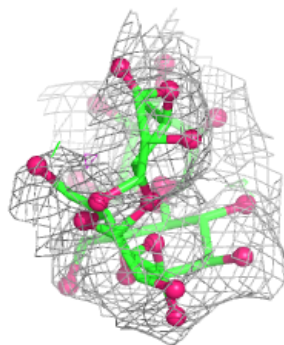
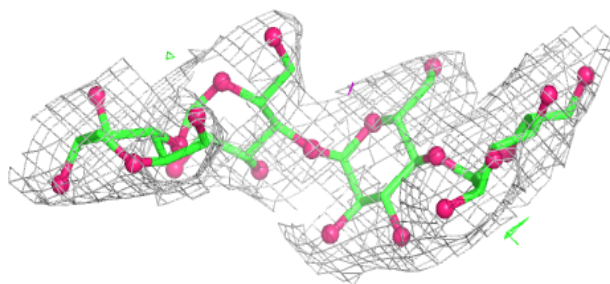
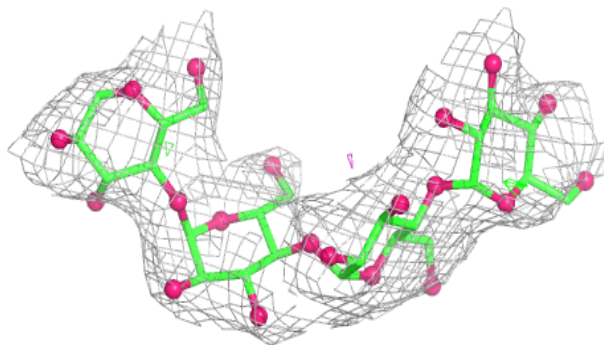
Electron density around Chain C:

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



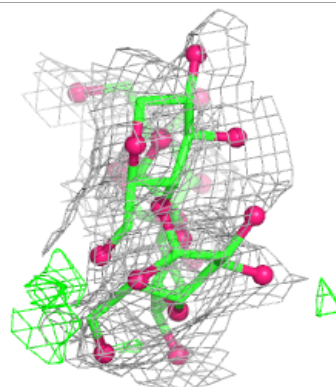
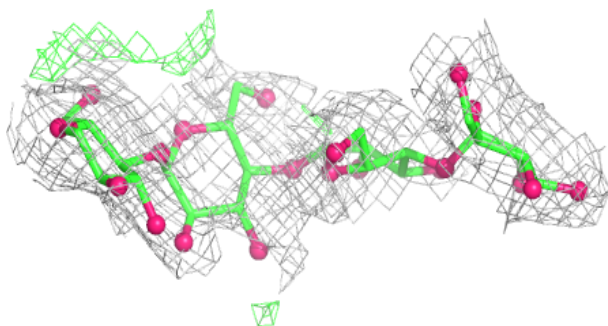
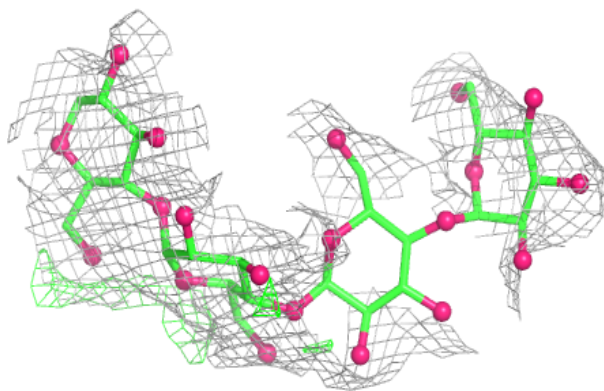
Electron density around Chain E:

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

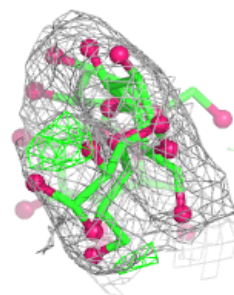
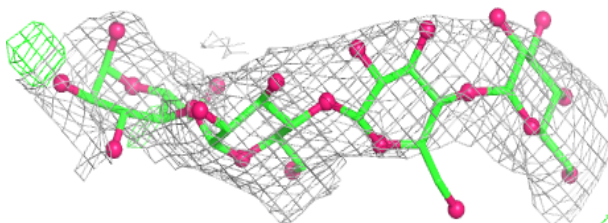
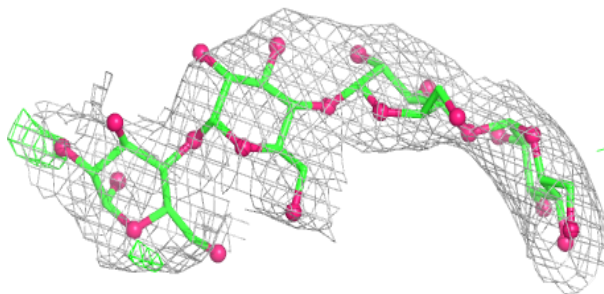


Electron density around Chain H:

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

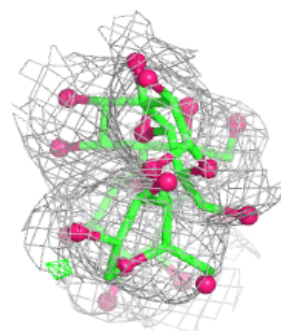
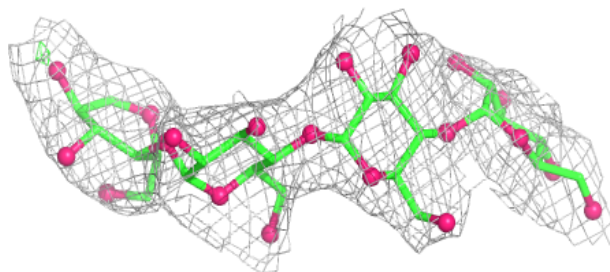
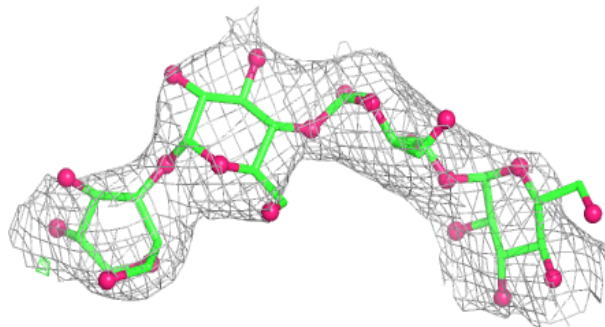
**Electron density around Chain I:**

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

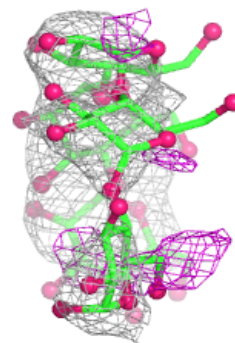
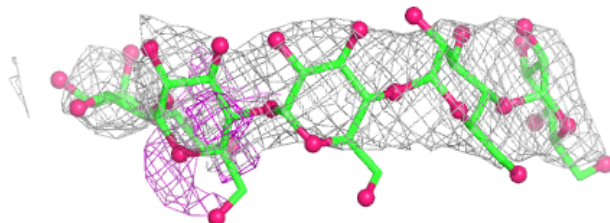
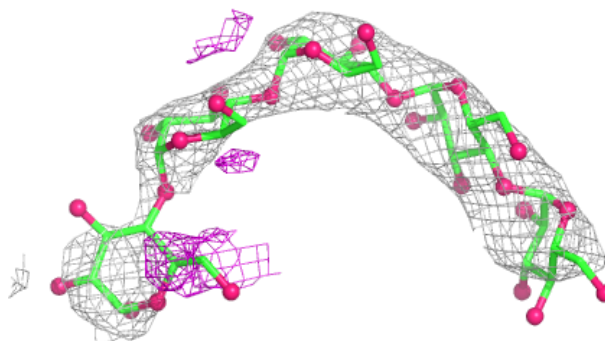


Electron density around Chain K:

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

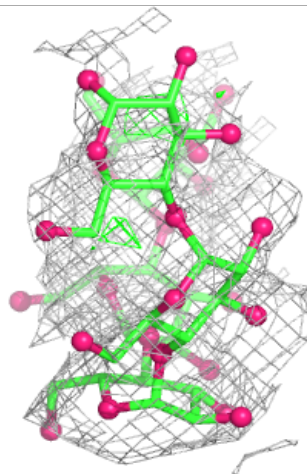
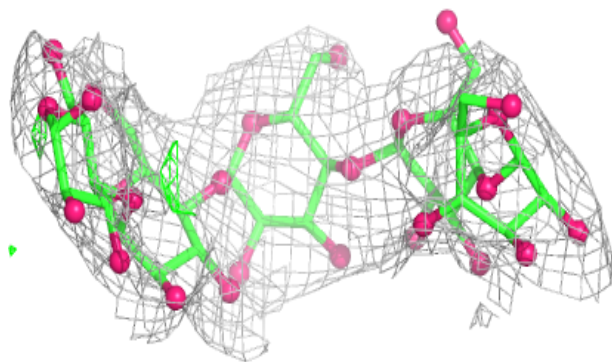
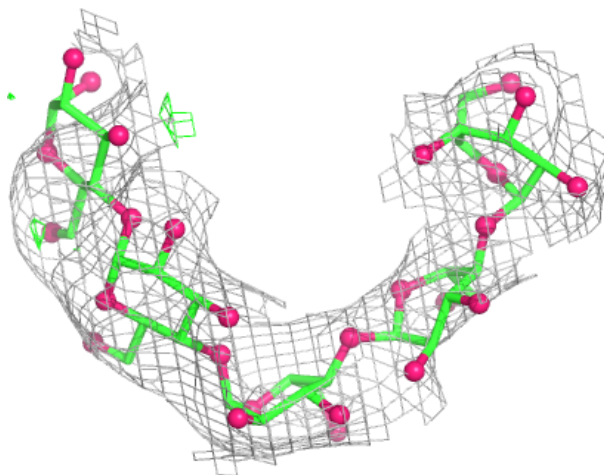
**Electron density around Chain D:**

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



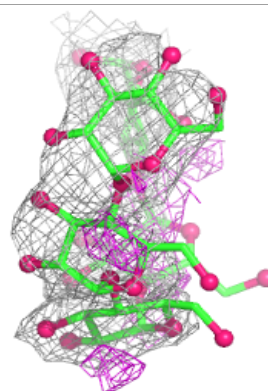
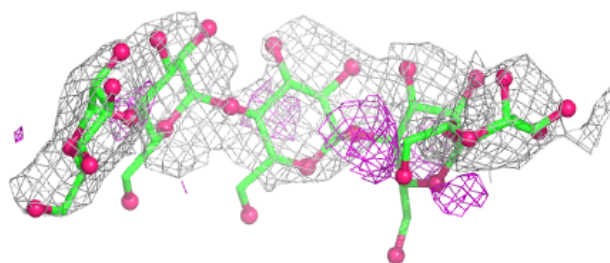
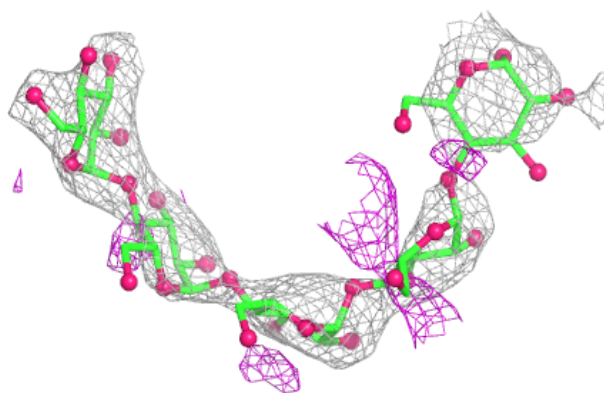
Electron density around Chain G:

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

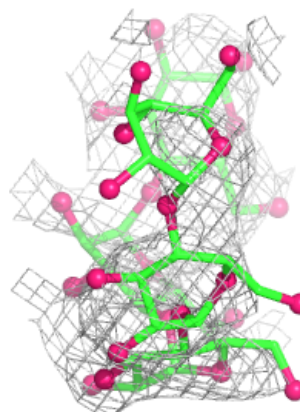
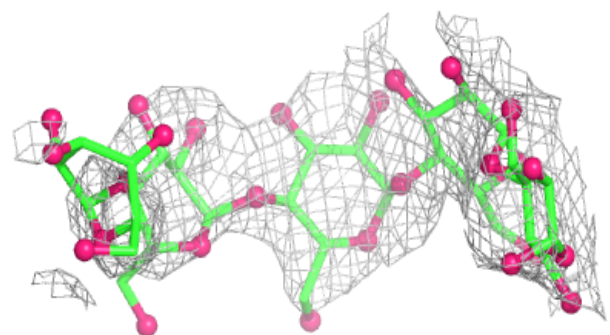
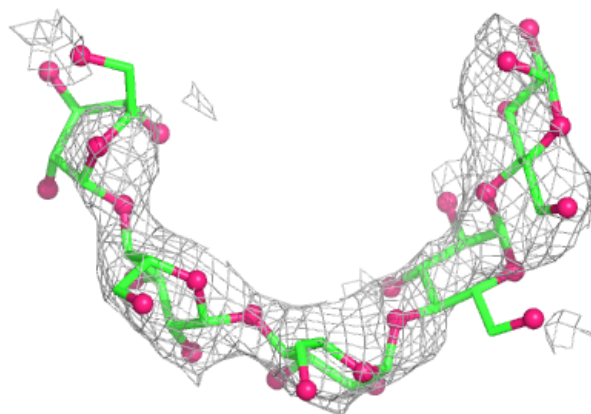


Electron density around Chain J:

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

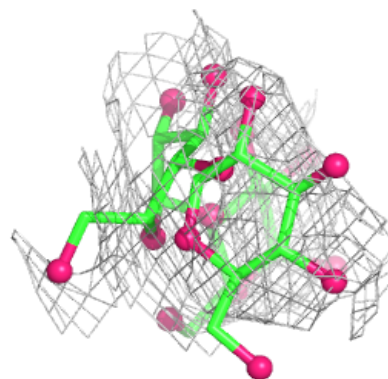
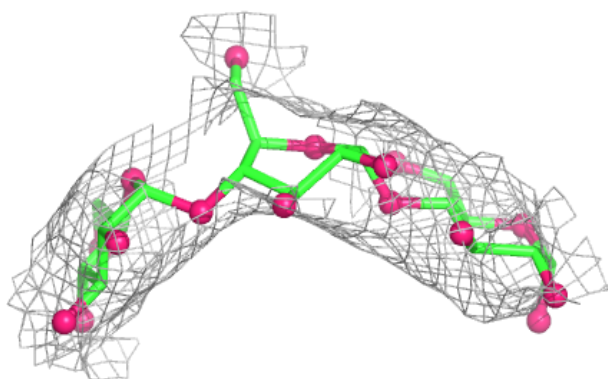
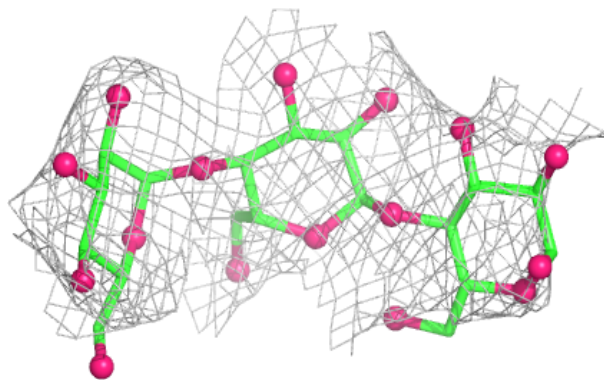
**Electron density around Chain M:**

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

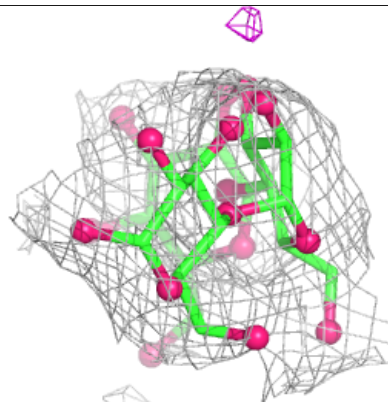
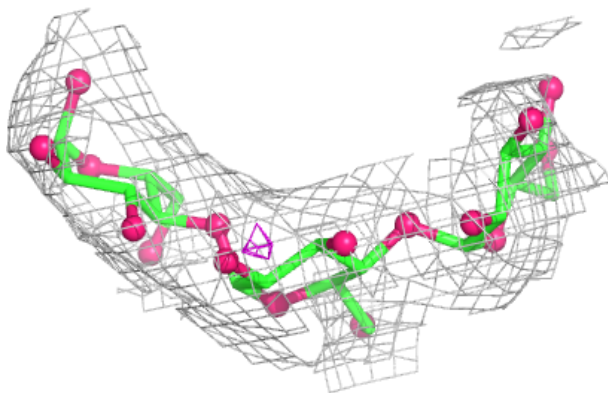
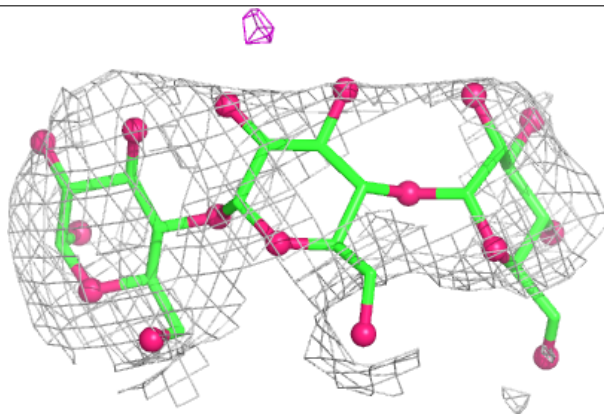


Electron density around Chain F:

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

**Electron density around Chain L:**

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



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.