



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 5, 2023 – 01:59 AM EST

PDB ID : 5I0D
Title : Cycloalternan-forming enzyme from *Listeria monocytogenes* in complex with cycloalternan
Authors : Light, S.H.; Minasov, G.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2016-02-03
Resolution : 1.77 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 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 : 4.02b-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

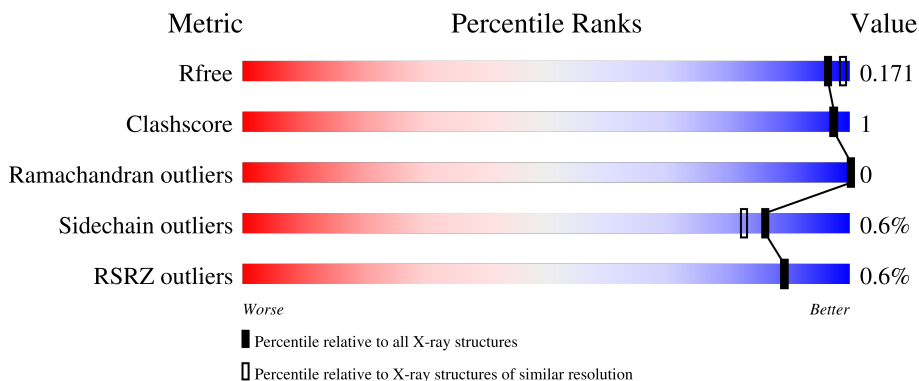
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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	1063	97%
1	B	1063	96%
2	C	2	100%
2	E	2	100%
2	I	2	100%

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Mol	Chain	Length	Quality of chain
2	J	2	 100%
3	D	4	 25% 75%
3	F	4	 25% 75%
3	K	4	 50% 50%
3	L	4	 75% 25%
3	M	4	 100%
4	G	3	 33% 67%
5	H	4	 50% 50%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 20432 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 Lmo2446 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	1060	8555	5365	1365	1791	34	0	29	0
1	B	1060	8610	5391	1370	1815	34	0	35	0

There are 4 discrepancies between the modelled and reference sequences:

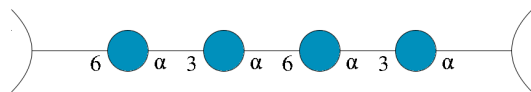
Chain	Residue	Modelled	Actual	Comment	Reference
A	29	SER	-	expression tag	UNP Q8Y4J2
A	30	ASN	-	expression tag	UNP Q8Y4J2
B	29	SER	-	expression tag	UNP Q8Y4J2
B	30	ASN	-	expression tag	UNP Q8Y4J2

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	C	2	23	12	11	0	0	0
2	E	2	23	12	11	0	0	0
2	I	2	23	12	11	0	0	0
2	J	2	23	12	11	0	0	0

- Molecule 3 is an oligosaccharide called Cyclic alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose.



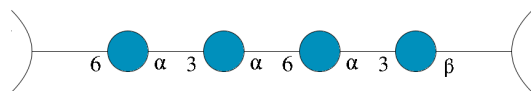
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	D	4	Total	C	O	0	0	0
			44	24	20			
3	F	4	Total	C	O	0	0	0
			44	24	20			
3	K	4	Total	C	O	0	0	0
			44	24	20			
3	L	4	Total	C	O	0	0	0
			44	24	20			
3	M	4	Total	C	O	0	0	0
			44	24	20			

- Molecule 4 is an oligosaccharide called alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose.



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

- Molecule 5 is an oligosaccharide called Cyclic alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
5	H	4	Total	C	O	0	0	0
			44	24	20			

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	5	Total	Mg	0	0
			5	5		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	3	Total	Mg	0	0
			3	3		

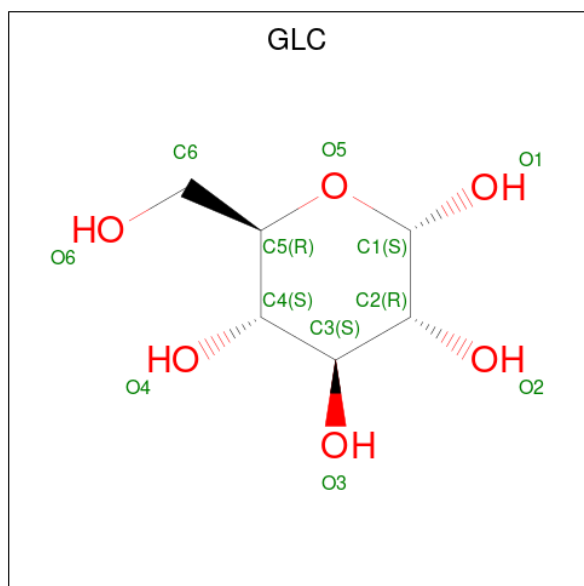
- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Ca	0	0
			1	1		
7	B	1	Total	Ca	0	0
			1	1		

- Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	2	Total	Cl	0	0
			2	2		
8	B	5	Total	Cl	0	0
			5	5		

- Molecule 9 is alpha-D-glucopyranose (three-letter code: GLC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			12	6	6		
9	B	1	Total	C	O	0	0
			12	6	6		

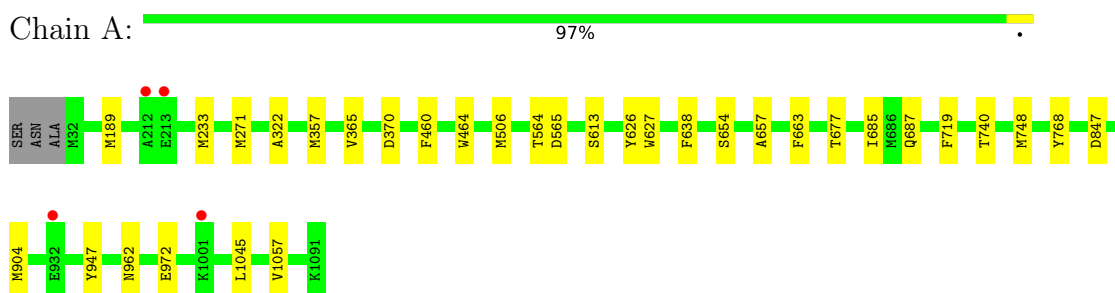
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	1415	Total 1424	O 1424	0	11
10	B	1398	Total 1412	O 1412	0	16

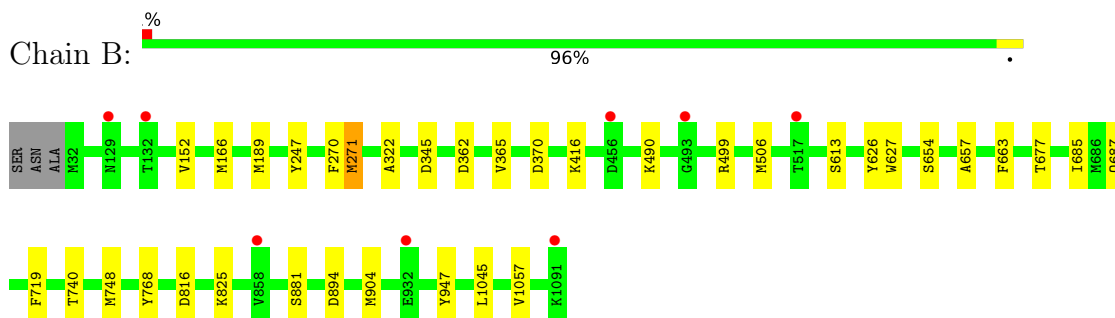
3 Residue-property plots [i](#)

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.

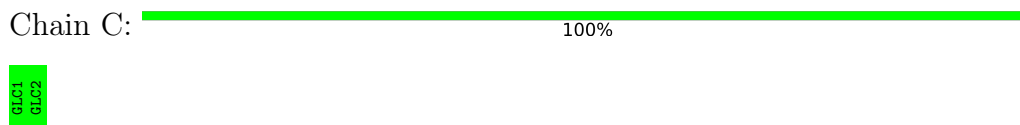
- Molecule 1: Lmo2446 protein



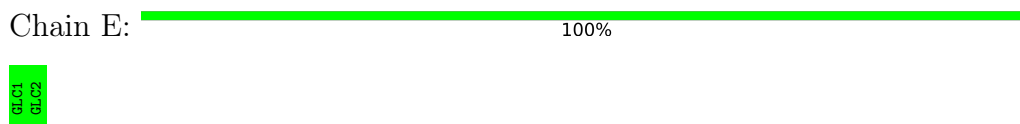
- Molecule 1: Lmo2446 protein



- Molecule 2: alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose



- Molecule 2: alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose



- Molecule 2: alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose



GLC1
GLC2

- Molecule 2: alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose

Chain J:  100%

GLC1
GLC2

- Molecule 3: Cyclic alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose

Chain D:  25% 75%

GLC1
GLC2
GLC3
GLC4

- Molecule 3: Cyclic alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose

Chain F:  25% 75%

GLC1
GLC2
GLC3
GLC4

- Molecule 3: Cyclic alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose

Chain K:  50% 50%

GLC1
GLC2
GLC3
GLC4

- Molecule 3: Cyclic alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose

Chain L:  75% 25%


GLC1
GLC2
GLC3
GLC4

- Molecule 3: Cyclic alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose

Chain M:  100%

GLC1
GLC2
GLC3
GLC4

- Molecule 4: alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose

Chain G:  33% 67%

GLC1
GLC2
GLC3

- Molecule 5: Cyclic alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose-(1-3)-alpha-D-glucopyranose-(1-6)-alpha-D-glucopyranose

Chain H:  50% 50%

GLC1
GLC2
GLC3
GLC4

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	74.75Å 101.23Å 166.39Å 90.00° 101.02° 90.00°	Depositor
Resolution (Å)	30.00 – 1.77 29.91 – 1.77	Depositor EDS
% Data completeness (in resolution range)	98.4 (30.00-1.77) 98.4 (29.91-1.77)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 1.77Å)	Xtrriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.144 , 0.171 0.144 , 0.171	Depositor DCC
R_{free} test set	11555 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	16.5	Xtrriage
Anisotropy	0.448	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.015 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	20432	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.15% 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: GLC, CA, MG, 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	0.50	0/8735	0.72	7/11846 (0.1%)
1	B	0.51	2/8790 (0.0%)	0.73	5/11924 (0.0%)
All	All	0.50	2/17525 (0.0%)	0.73	12/23770 (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	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	271[A]	MSE	N-CA	8.33	1.63	1.46
1	B	271[B]	MSE	N-CA	8.33	1.63	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	271[A]	MSE	CB-CG-SE	8.28	137.54	112.70
1	B	271[B]	MSE	CB-CG-SE	8.28	137.54	112.70
1	A	972	GLU	OE1-CD-OE2	6.77	131.42	123.30
1	A	565	ASP	CB-CG-OD1	6.75	124.38	118.30
1	A	370	ASP	CB-CG-OD1	6.46	124.11	118.30
1	A	847	ASP	CB-CG-OD1	5.58	123.32	118.30
1	A	638	PHE	CB-CG-CD1	5.45	124.62	120.80
1	A	638	PHE	CB-CG-CD2	-5.38	117.03	120.80
1	A	357	MSE	CG-SE-CE	-5.38	87.07	98.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	816	ASP	CB-CG-OD1	5.20	122.98	118.30
1	B	362	ASP	CB-CG-OD1	5.08	122.87	118.30
1	B	370	ASP	CB-CG-OD1	5.06	122.85	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	270	PHE	Mainchain

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	8555	0	7851	13	0
1	B	8610	0	7863	15	0
2	C	23	0	21	0	0
2	E	23	0	21	0	0
2	I	23	0	21	0	0
2	J	23	0	21	0	0
3	D	44	0	36	0	0
3	F	44	0	36	0	0
3	K	44	0	36	0	0
3	L	44	0	36	0	0
3	M	44	0	36	0	0
4	G	34	0	30	0	0
5	H	44	0	36	0	0
6	A	5	0	0	0	0
6	B	3	0	0	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
8	A	2	0	0	0	0
8	B	5	0	0	0	0
9	A	12	0	12	0	0
9	B	12	0	12	0	0
10	A	1424	0	0	3	1
10	B	1412	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	20432	0	16068	28	1

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 1.

All (28) 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:189[B]:MSE:HE1	1:A:365:VAL:HG23	1.70	0.73
1:B:271[B]:MSE:HA	1:B:345:ASP:O	2.08	0.54
1:B:613:SER:O	1:B:626:TYR:HA	2.10	0.52
1:B:748:MSE:HG3	1:B:768:TYR:CD1	2.45	0.51
1:B:657:ALA:HB2	1:B:685:ILE:HB	1.93	0.50
1:B:1045:LEU:HD23	1:B:1057[A]:VAL:HG21	1.93	0.50
1:B:152:VAL:HG22	1:B:166[A]:MSE:SE	2.62	0.50
1:A:271[B]:MSE:HE2	10:A:1645:HOH:O	2.12	0.50
1:B:189[B]:MSE:HE1	1:B:365:VAL:HG23	1.94	0.50
1:A:613:SER:O	1:A:626:TYR:HA	2.11	0.50
1:A:657:ALA:HB2	1:A:685:ILE:HB	1.95	0.48
1:B:904:MSE:O	1:B:947:TYR:HA	2.13	0.48
1:B:271[A]:MSE:HE2	10:B:1963:HOH:O	2.14	0.48
1:B:654:SER:HB3	1:B:740:THR:HG21	1.96	0.47
1:B:881[A]:SER:OG	1:B:894[A]:ASP:HB2	2.16	0.46
1:A:654:SER:HB3	1:A:740:THR:HG21	1.98	0.46
1:B:490:LYS:O	1:B:499:ARG:HD3	2.17	0.45
1:A:748:MSE:HG3	1:A:768:TYR:CD1	2.51	0.45
1:A:460:PHE:HB3	1:A:464:TRP:O	2.18	0.44
1:A:677:THR:HG21	1:A:719:PHE:HA	1.99	0.43
1:B:247:TYR:CE1	1:B:271[B]:MSE:HB2	2.54	0.43
1:A:322:ALA:HB3	10:A:1520:HOH:O	2.19	0.43
1:B:677:THR:HG21	1:B:719:PHE:HA	2.02	0.42
1:B:322:ALA:HB3	10:B:2196:HOH:O	2.18	0.42
1:A:904:MSE:O	1:A:947:TYR:HA	2.20	0.41
1:A:962:ASN:ND2	10:A:1222:HOH:O	2.54	0.41
1:A:189[B]:MSE:SE	1:A:233:MSE:SE	3.40	0.40
1:A:1045:LEU:HD23	1:A:1057[A]:VAL:HG21	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:A:2374:HOH:O	10:A:2374:HOH:O[2_857]	2.17	0.03

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	1086/1063 (102%)	1064 (98%)	22 (2%)	0	100	100
1	B	1093/1063 (103%)	1063 (97%)	30 (3%)	0	100	100
All	All	2179/2126 (102%)	2127 (98%)	52 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	919/863 (106%)	914 (100%)	5 (0%)	88	86
1	B	926/863 (107%)	920 (99%)	6 (1%)	86	82
All	All	1845/1726 (107%)	1834 (99%)	11 (1%)	86	82

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	506	MSE
1	A	564	THR
1	A	627	TRP

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Mol	Chain	Res	Type
1	A	663	PHE
1	A	687	GLN
1	B	416	LYS
1	B	506	MSE
1	B	627	TRP
1	B	663	PHE
1	B	687	GLN
1	B	825	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	304	GLN
1	A	965	ASN
1	A	975	HIS
1	A	1064	GLN
1	B	965	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](#)

35 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	0.61	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GLC	C	2	2	11,11,12	0.18	0	15,15,17	0.72	0
3	GLC	D	1	3	11,11,12	0.31	0	15,15,17	0.95	1 (6%)
3	GLC	D	2	3	11,11,12	0.31	0	15,15,17	0.80	1 (6%)
3	GLC	D	3	3	11,11,12	0.42	0	15,15,17	1.05	2 (13%)
3	GLC	D	4	3	11,11,12	0.20	0	15,15,17	0.84	0
2	GLC	E	1	2	12,12,12	0.51	0	17,17,17	0.62	0
2	GLC	E	2	2	11,11,12	0.29	0	15,15,17	0.70	0
3	GLC	F	1	3	11,11,12	0.40	0	15,15,17	0.97	1 (6%)
3	GLC	F	2	3	11,11,12	0.41	0	15,15,17	0.91	1 (6%)
3	GLC	F	3	3	11,11,12	0.68	0	15,15,17	1.18	0
3	GLC	F	4	3	11,11,12	0.44	0	15,15,17	1.17	2 (13%)
4	GLC	G	1	4	12,12,12	0.51	0	17,17,17	0.61	0
4	GLC	G	2	4	11,11,12	0.38	0	15,15,17	1.14	1 (6%)
4	GLC	G	3	4	11,11,12	0.21	0	15,15,17	0.86	1 (6%)
5	GLC	H	1	5	11,11,12	0.52	0	15,15,17	0.93	1 (6%)
5	GLC	H	2	5	11,11,12	0.38	0	15,15,17	0.93	0
5	GLC	H	3	5	11,11,12	0.58	0	15,15,17	0.84	1 (6%)
5	GLC	H	4	5	11,11,12	0.67	0	15,15,17	0.96	0
2	GLC	I	1	2	12,12,12	0.46	0	17,17,17	0.48	0
2	GLC	I	2	2	11,11,12	0.29	0	15,15,17	0.74	0
2	GLC	J	1	2	12,12,12	0.52	0	17,17,17	0.64	0
2	GLC	J	2	2	11,11,12	0.31	0	15,15,17	0.81	0
3	GLC	K	1	3	11,11,12	0.44	0	15,15,17	1.10	1 (6%)
3	GLC	K	2	3	11,11,12	0.32	0	15,15,17	0.89	1 (6%)
3	GLC	K	3	3	11,11,12	0.54	0	15,15,17	1.01	0
3	GLC	K	4	3	11,11,12	0.46	0	15,15,17	0.72	0
3	GLC	L	1	3	11,11,12	0.46	0	15,15,17	0.86	0
3	GLC	L	2	3	11,11,12	0.54	0	15,15,17	0.95	0
3	GLC	L	3	3	11,11,12	0.54	0	15,15,17	0.86	0
3	GLC	L	4	3	11,11,12	0.46	0	15,15,17	0.87	1 (6%)
3	GLC	M	1	3	11,11,12	0.37	0	15,15,17	1.09	2 (13%)
3	GLC	M	2	3	11,11,12	0.27	0	15,15,17	0.99	1 (6%)
3	GLC	M	3	3	11,11,12	0.38	0	15,15,17	0.98	1 (6%)
3	GLC	M	4	3	11,11,12	0.26	0	15,15,17	0.75	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. '2' 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	GLC	C	2	2	-	0/2/19/22	0/1/1/1
3	GLC	D	1	3	-	0/2/19/22	0/1/1/1
3	GLC	D	2	3	-	0/2/19/22	0/1/1/1
3	GLC	D	3	3	-	0/2/19/22	0/1/1/1
3	GLC	D	4	3	-	0/2/19/22	0/1/1/1
2	GLC	E	1	2	-	0/2/22/22	0/1/1/1
2	GLC	E	2	2	-	0/2/19/22	0/1/1/1
3	GLC	F	1	3	-	0/2/19/22	0/1/1/1
3	GLC	F	2	3	-	0/2/19/22	0/1/1/1
3	GLC	F	3	3	-	0/2/19/22	0/1/1/1
3	GLC	F	4	3	-	0/2/19/22	0/1/1/1
4	GLC	G	1	4	-	0/2/22/22	0/1/1/1
4	GLC	G	2	4	-	1/2/19/22	0/1/1/1
4	GLC	G	3	4	-	0/2/19/22	0/1/1/1
5	GLC	H	1	5	-	0/2/19/22	0/1/1/1
5	GLC	H	2	5	-	0/2/19/22	0/1/1/1
5	GLC	H	3	5	-	0/2/19/22	0/1/1/1
5	GLC	H	4	5	-	0/2/19/22	0/1/1/1
2	GLC	I	1	2	-	0/2/22/22	0/1/1/1
2	GLC	I	2	2	-	0/2/19/22	0/1/1/1
2	GLC	J	1	2	-	0/2/22/22	0/1/1/1
2	GLC	J	2	2	-	0/2/19/22	0/1/1/1
3	GLC	K	1	3	-	1/2/19/22	0/1/1/1
3	GLC	K	2	3	-	0/2/19/22	0/1/1/1
3	GLC	K	3	3	-	0/2/19/22	0/1/1/1
3	GLC	K	4	3	-	0/2/19/22	0/1/1/1
3	GLC	L	1	3	-	2/2/19/22	0/1/1/1
3	GLC	L	2	3	-	0/2/19/22	0/1/1/1
3	GLC	L	3	3	-	0/2/19/22	0/1/1/1
3	GLC	L	4	3	-	0/2/19/22	0/1/1/1
3	GLC	M	1	3	-	0/2/19/22	0/1/1/1
3	GLC	M	2	3	-	0/2/19/22	0/1/1/1
3	GLC	M	3	3	-	0/2/19/22	0/1/1/1
3	GLC	M	4	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	2	GLC	C1-O5-C5	3.43	116.84	112.19
3	K	1	GLC	C1-O5-C5	2.87	116.08	112.19
3	F	2	GLC	C1-O5-C5	2.75	115.92	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	M	3	GLC	C1-O5-C5	2.52	115.60	112.19
3	K	2	GLC	C1-O5-C5	2.51	115.59	112.19
3	F	1	GLC	C1-O5-C5	2.51	115.59	112.19
3	D	1	GLC	C1-O5-C5	2.50	115.58	112.19
3	F	4	GLC	C1-O5-C5	2.38	115.42	112.19
5	H	3	GLC	C1-O5-C5	2.34	115.36	112.19
3	D	2	GLC	C1-O5-C5	2.29	115.30	112.19
3	D	3	GLC	C1-O5-C5	2.21	115.18	112.19
3	M	1	GLC	O5-C1-C2	-2.13	107.49	110.77
3	L	4	GLC	O5-C1-C2	-2.13	107.49	110.77
3	M	1	GLC	O3-C3-C2	-2.12	105.93	109.99
5	H	1	GLC	C1-O5-C5	2.08	115.01	112.19
3	D	3	GLC	O3-C3-C2	-2.08	106.02	109.99
3	M	2	GLC	C2-C3-C4	-2.06	107.32	110.89
4	G	3	GLC	C1-O5-C5	2.06	114.98	112.19
3	M	4	GLC	C1-O5-C5	2.05	114.97	112.19
3	F	4	GLC	O6-C6-C5	-2.03	104.33	111.29

There are no chirality outliers.

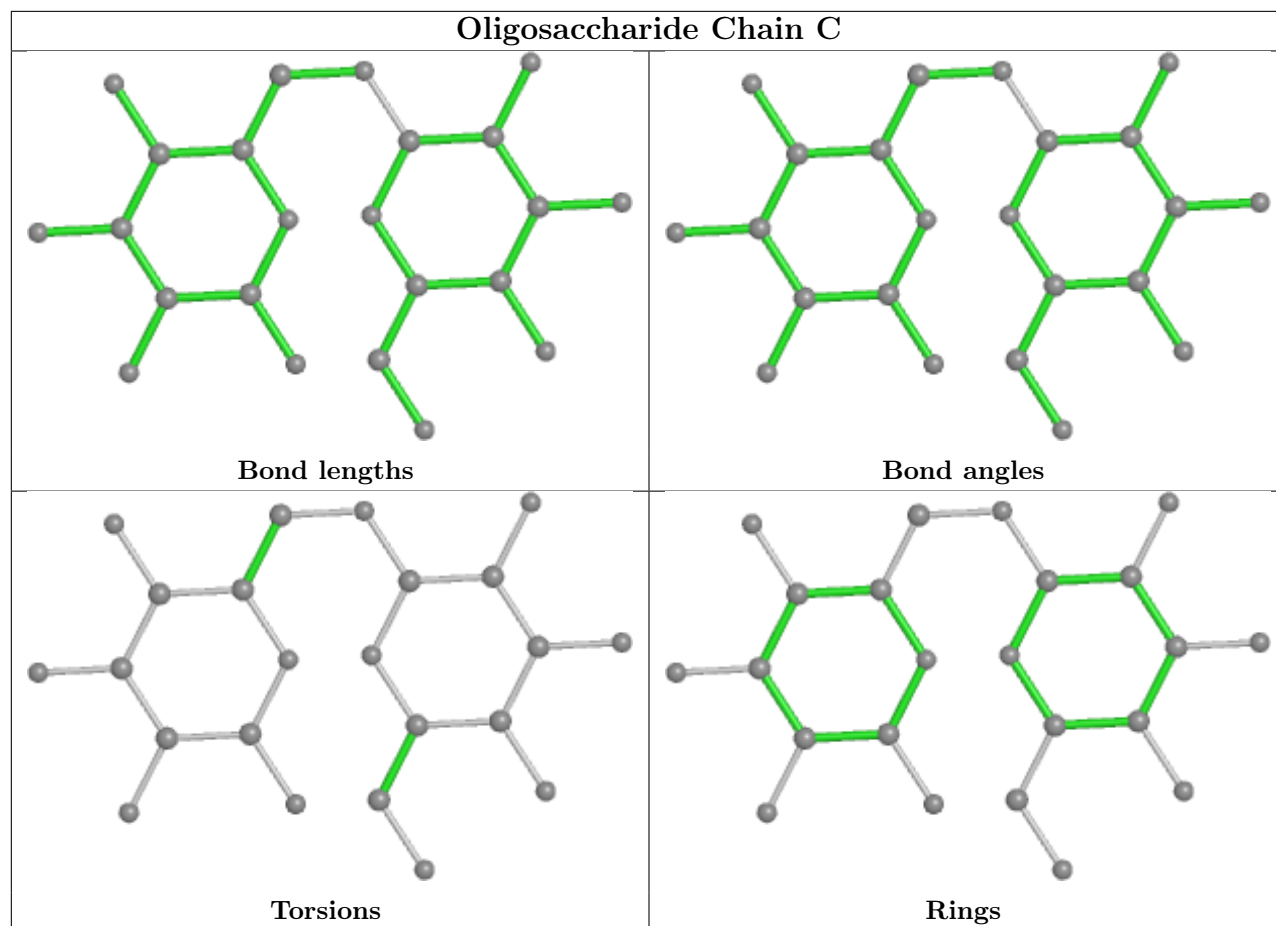
All (4) torsion outliers are listed below:

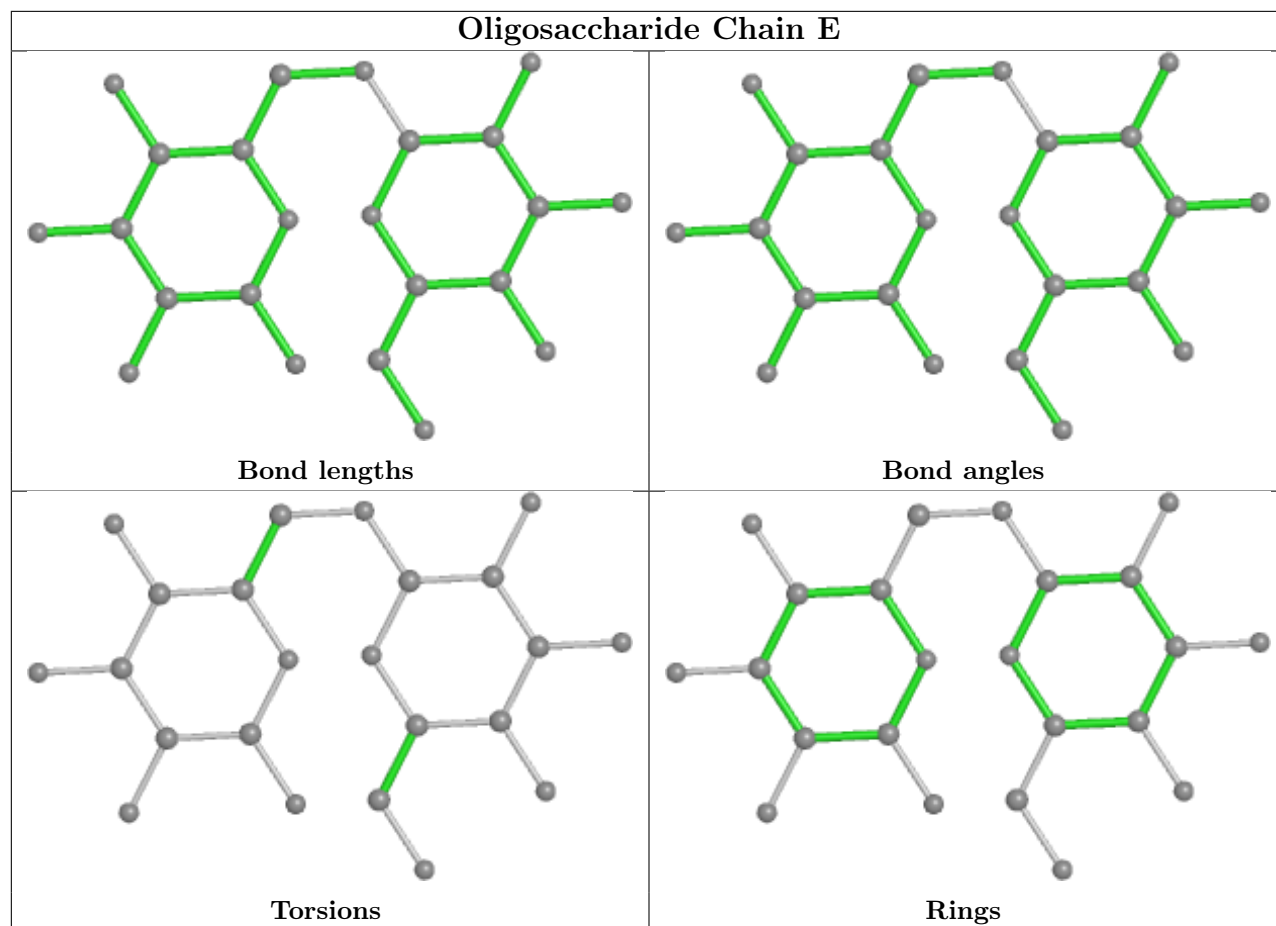
Mol	Chain	Res	Type	Atoms
3	L	1	GLC	C4-C5-C6-O6
3	L	1	GLC	O5-C5-C6-O6
3	K	1	GLC	C4-C5-C6-O6
4	G	2	GLC	C4-C5-C6-O6

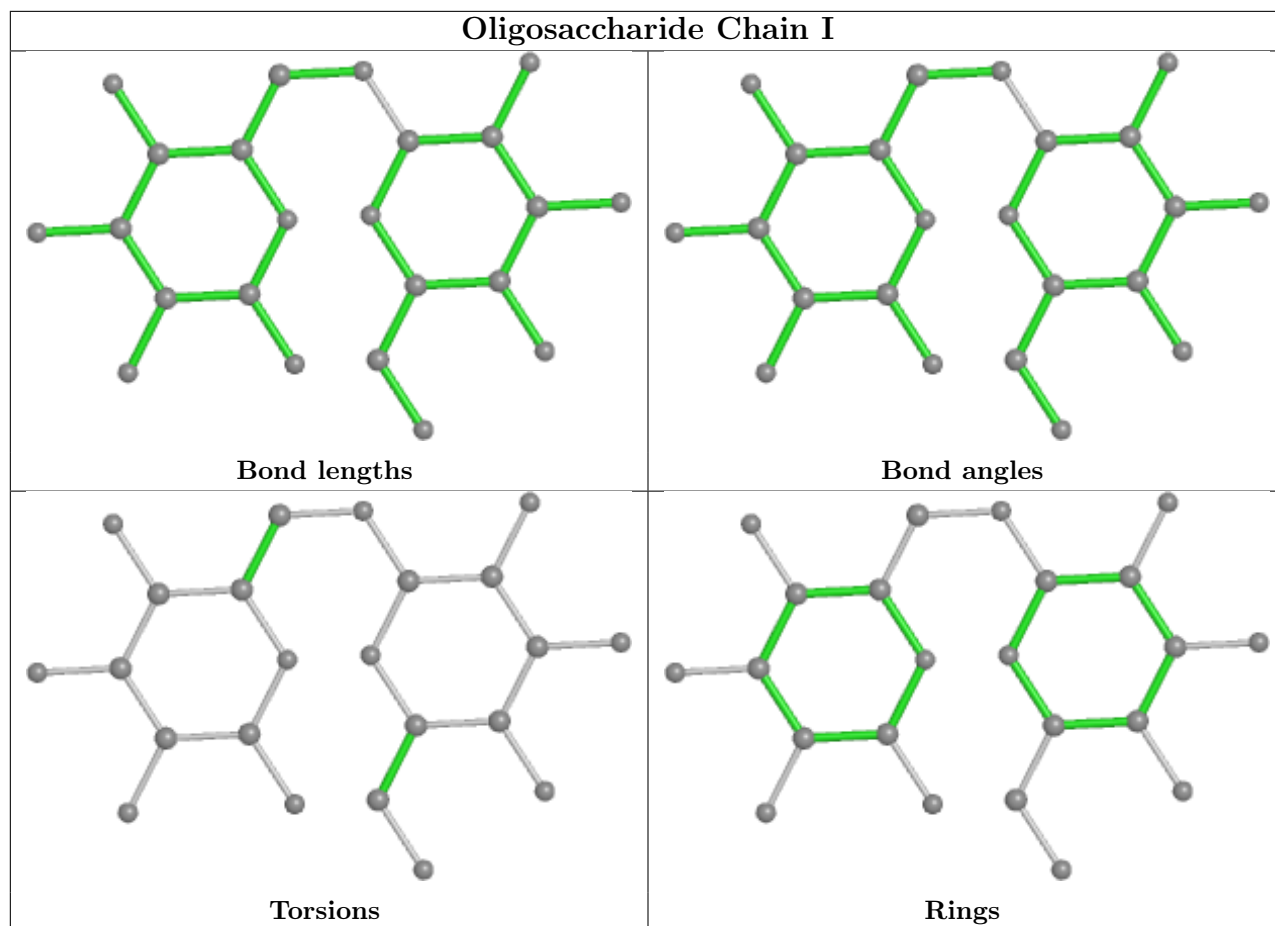
There are no ring outliers.

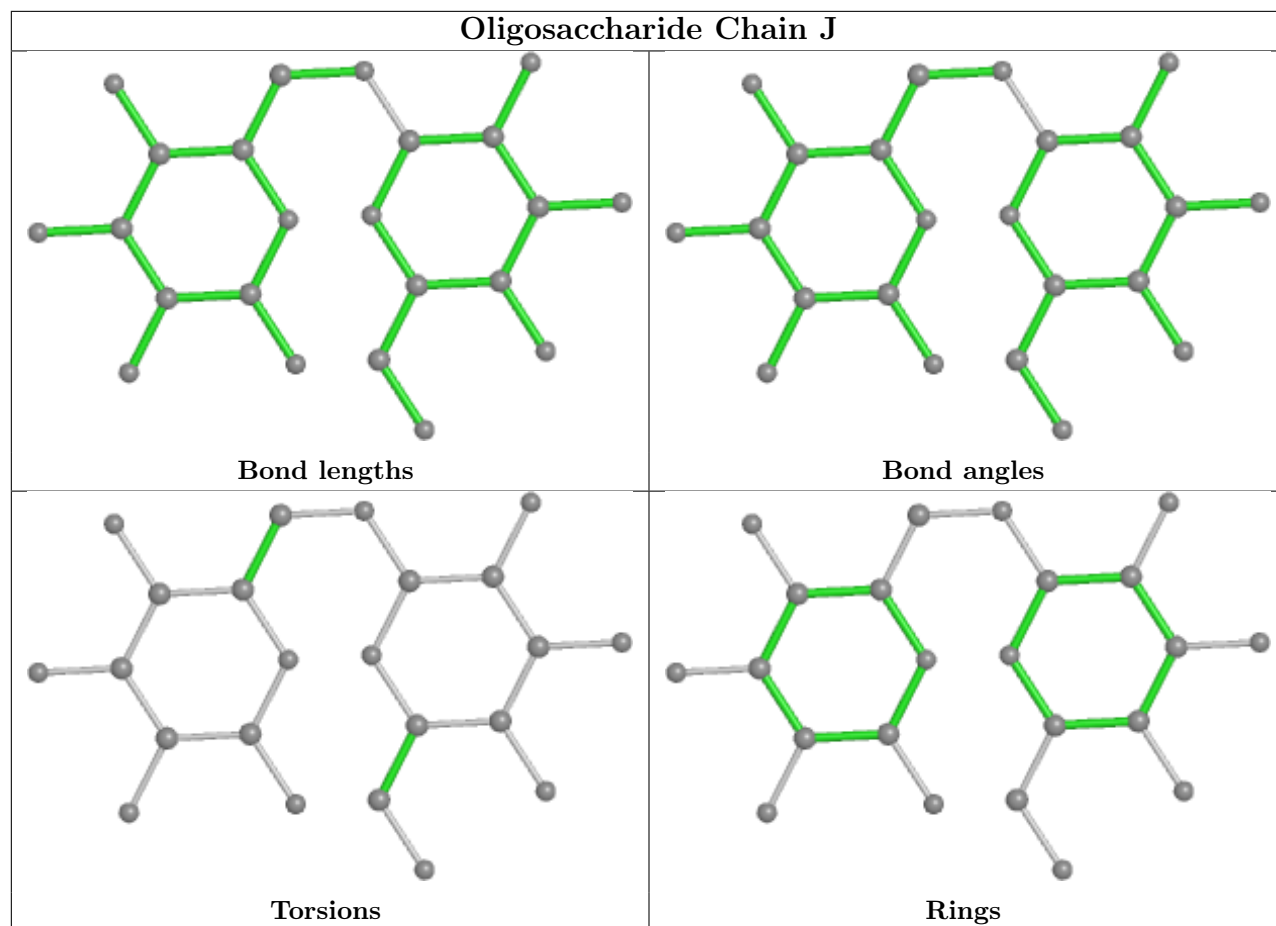
No monomer is involved in short contacts.

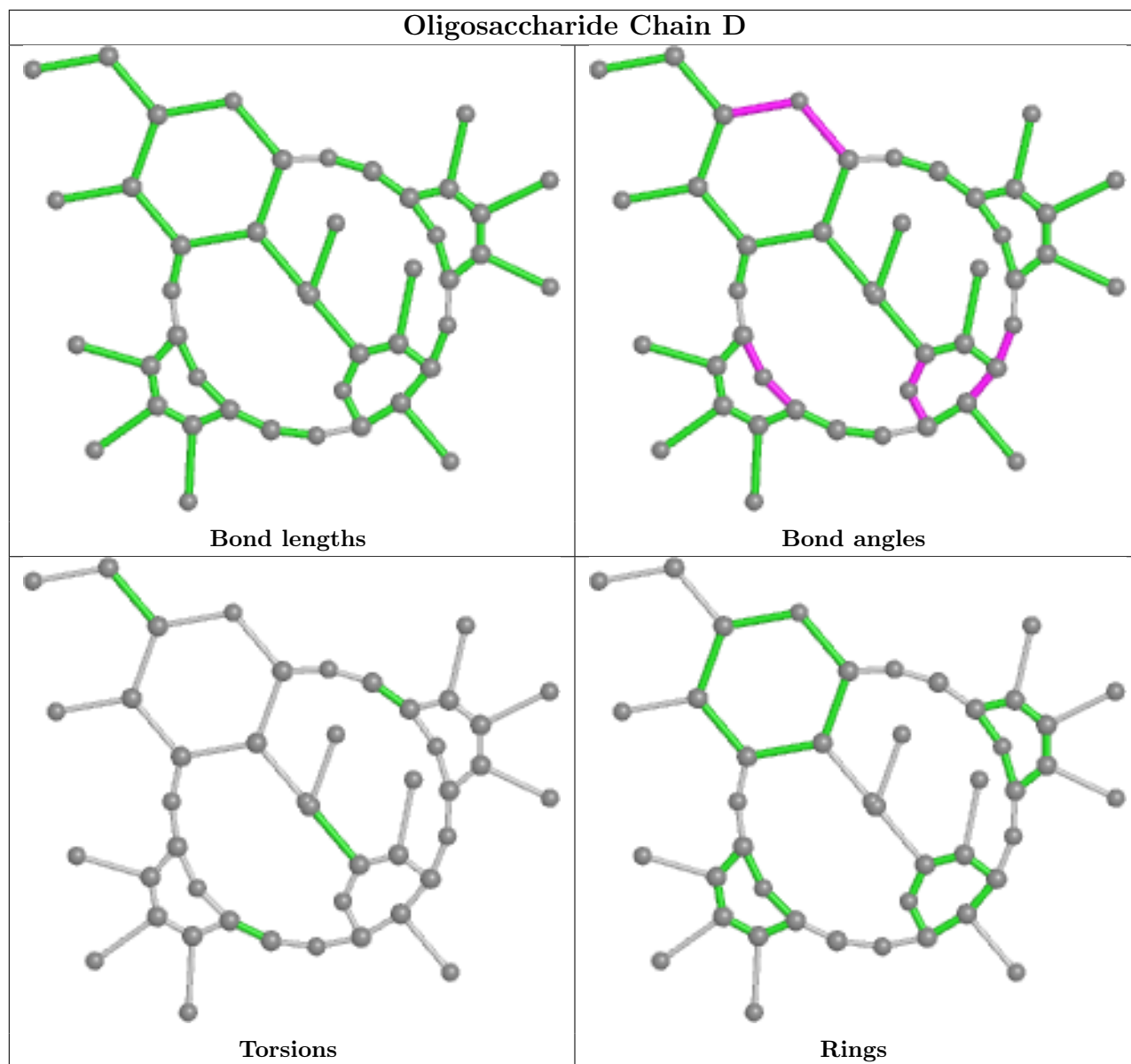
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

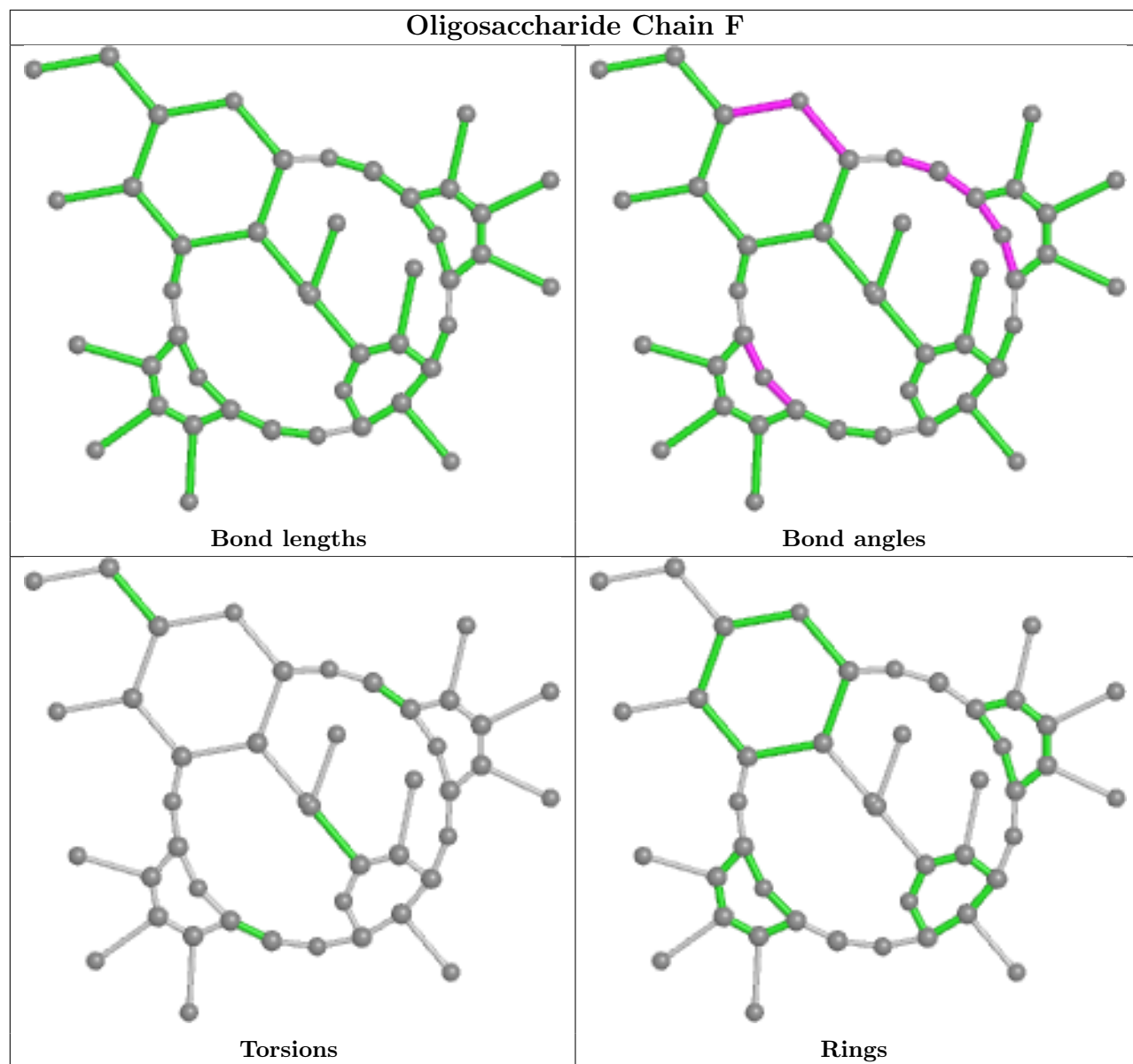


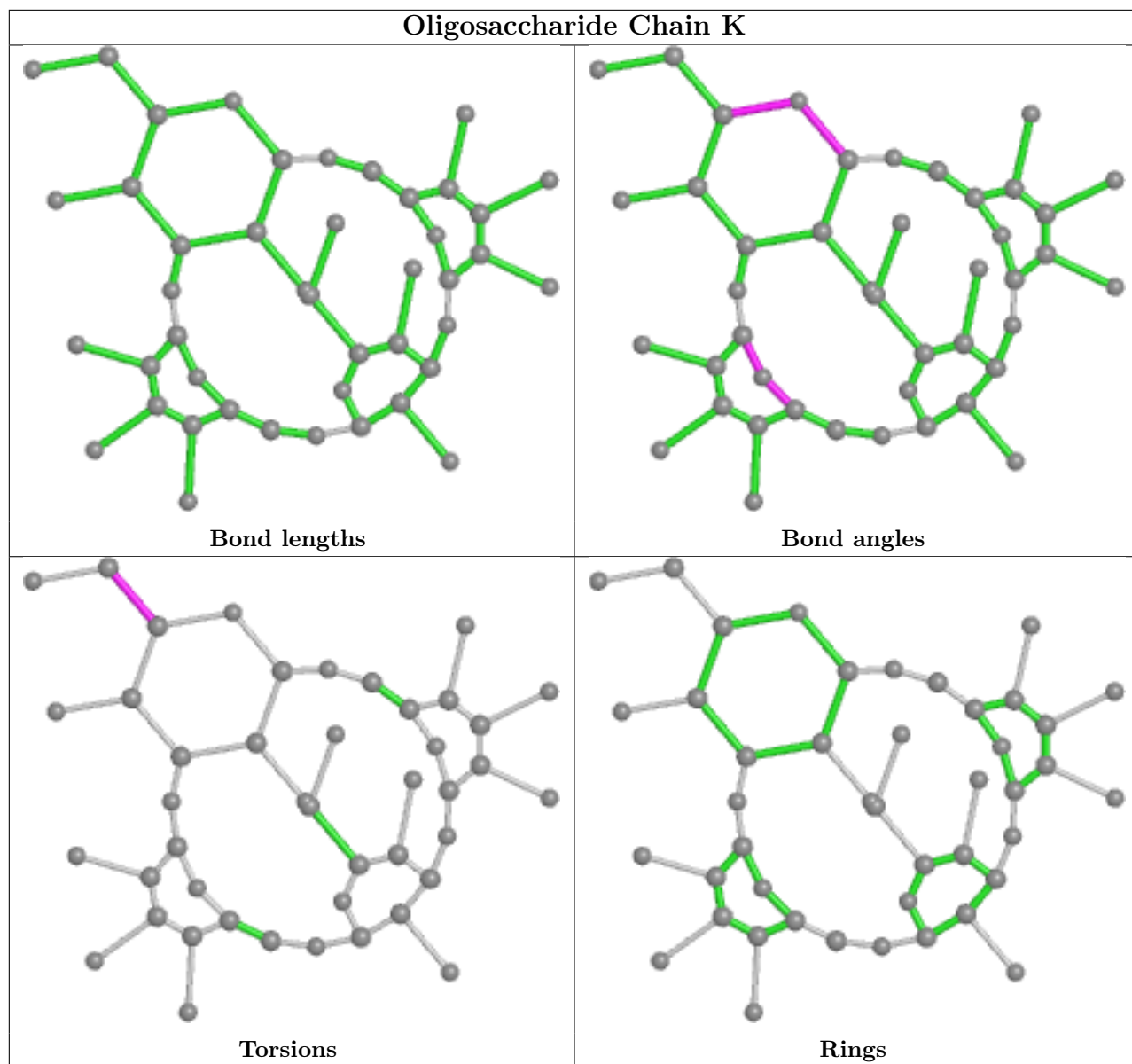


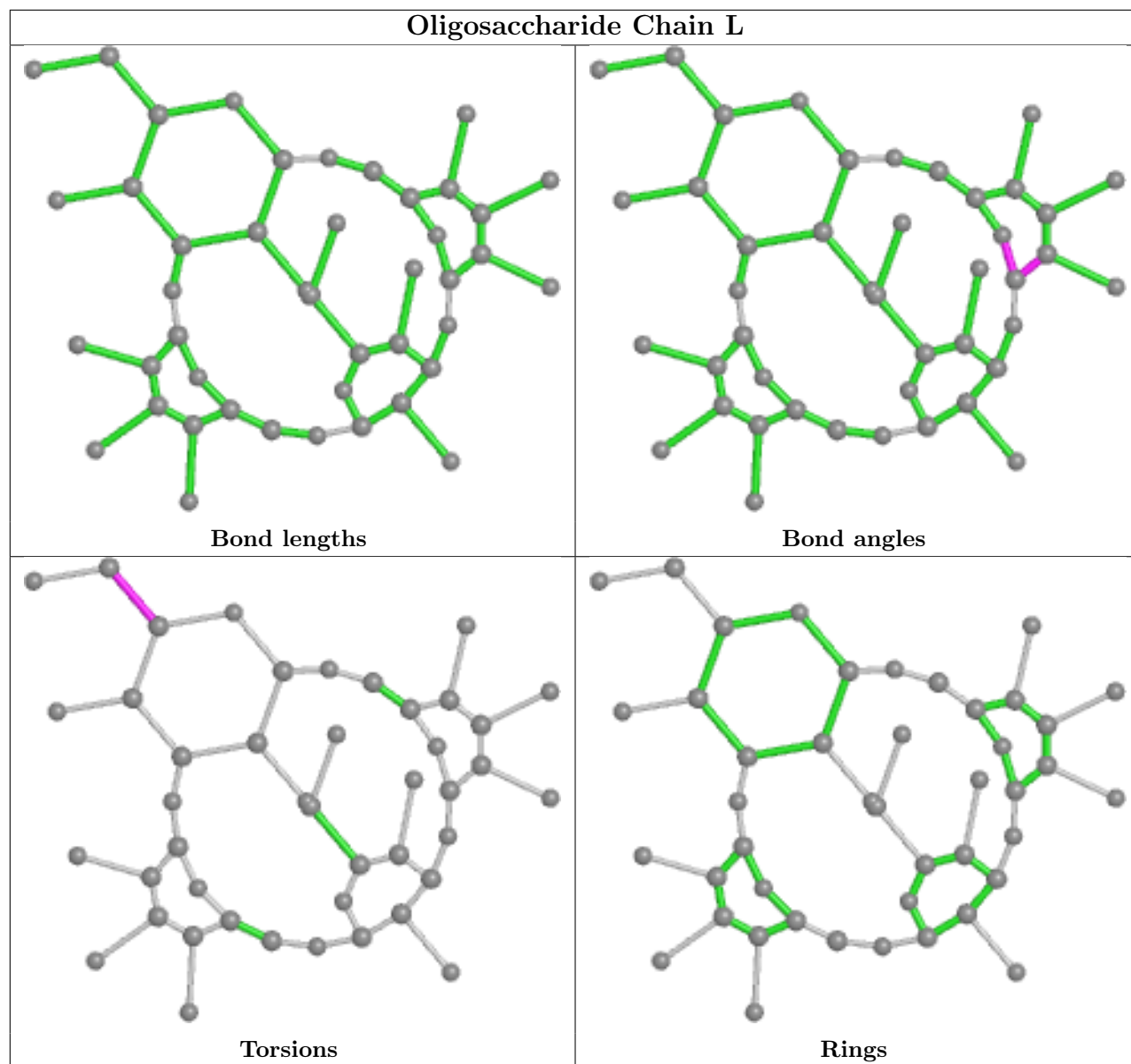


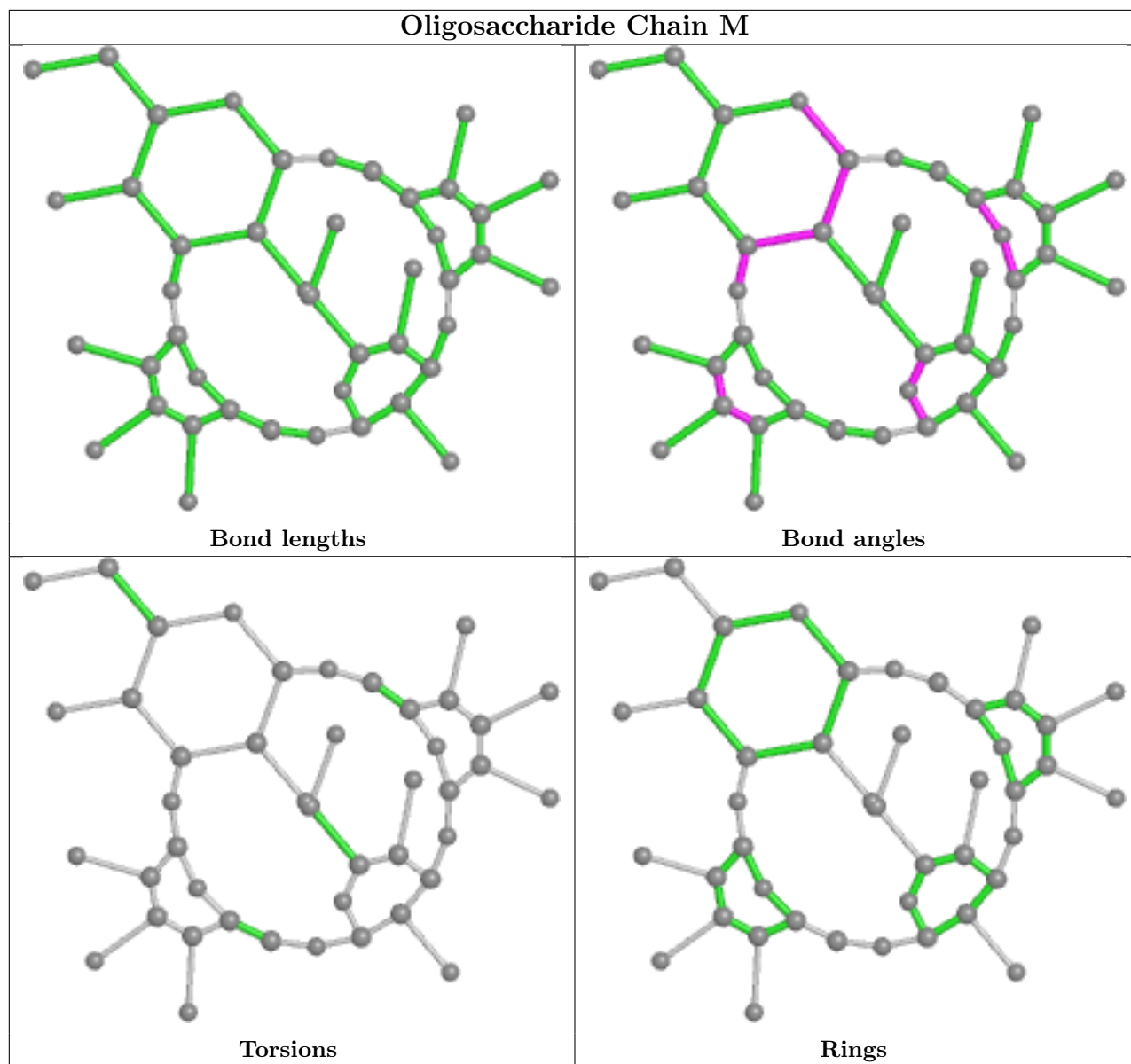


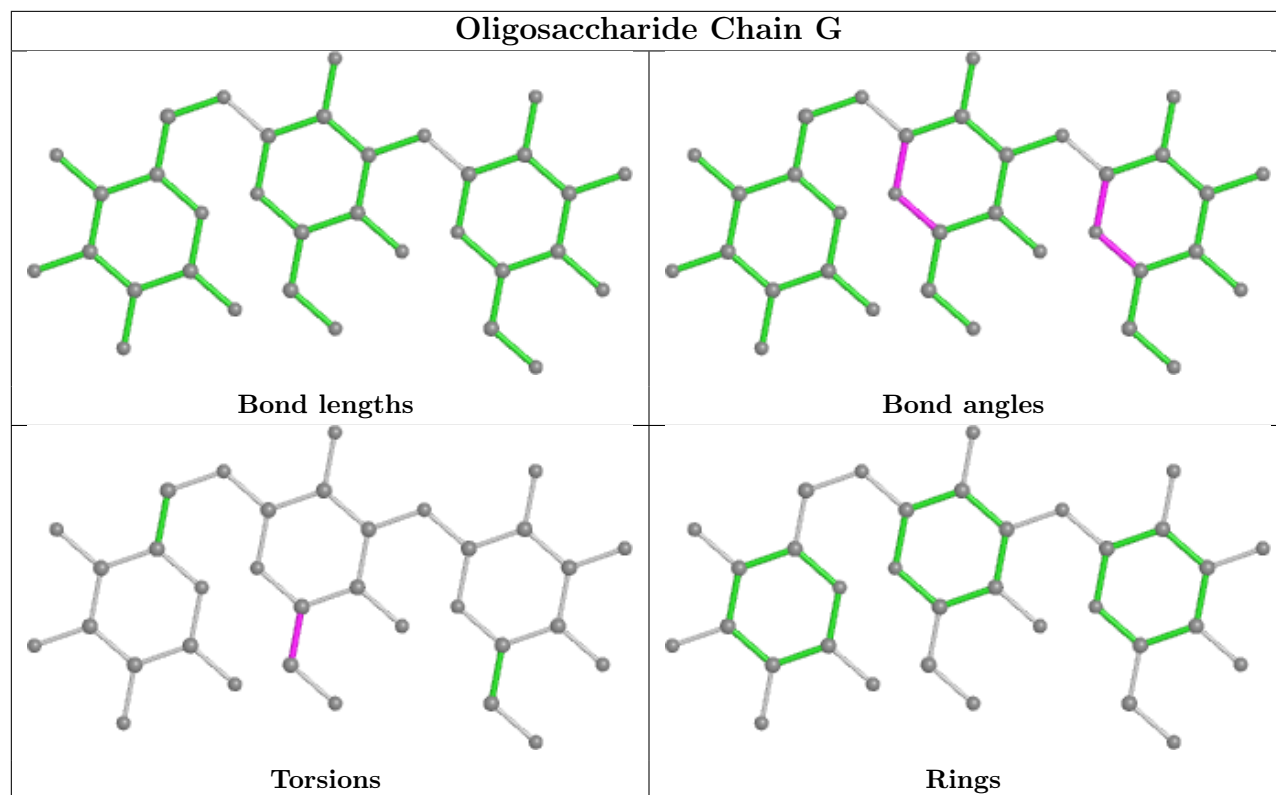


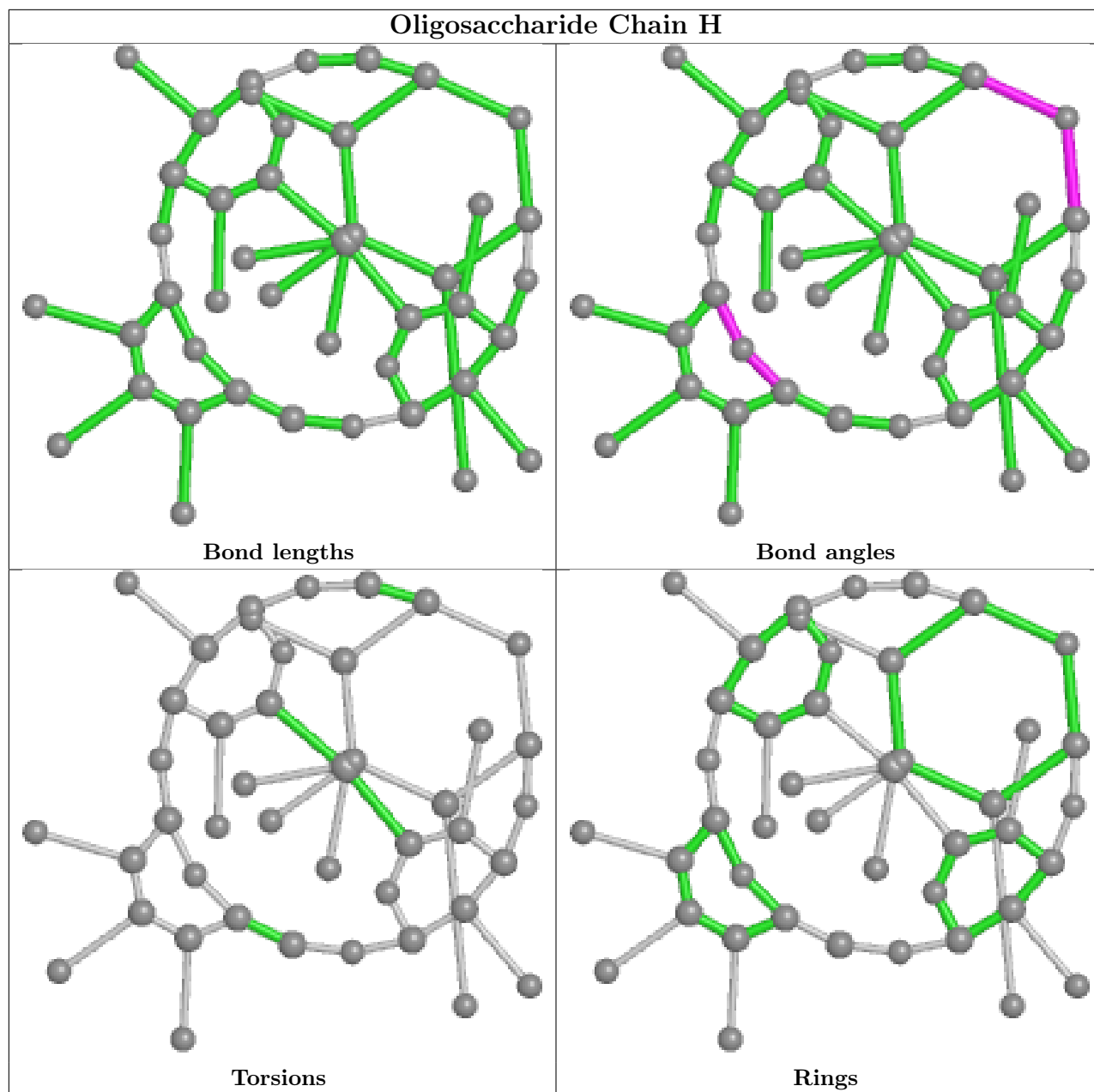












5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 17 are monoatomic - leaving 2 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
9	GLC	B	1122	-	12,12,12	0.49	0	17,17,17	0.56	0
9	GLC	A	1124	-	12,12,12	0.53	0	17,17,17	0.70	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
9	GLC	B	1122	-	-	0/2/22/22	0/1/1/1
9	GLC	A	1124	-	-	1/2/22/22	0/1/1/1

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
9	A	1124	GLC	C4-C5-C6-O6

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, 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	1030/1063 (96%)	-0.32	4 (0%) 92 92	11, 18, 30, 46	0
1	B	1030/1063 (96%)	-0.31	8 (0%) 86 86	11, 19, 30, 41	0
All	All	2060/2126 (96%)	-0.31	12 (0%) 89 89	11, 18, 30, 46	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	132	THR	4.6
1	A	212	ALA	3.5
1	B	932	GLU	3.1
1	B	129	ASN	3.0
1	B	858	VAL	2.8
1	B	493	GLY	2.5
1	A	932	GLU	2.5
1	B	1091	LYS	2.4
1	B	517	THR	2.1
1	A	1001	LYS	2.1
1	B	456[A]	ASP	2.1
1	A	213	GLU	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](#)

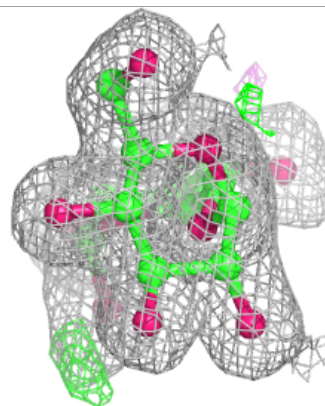
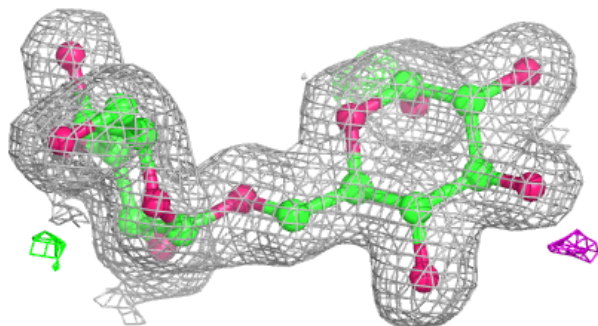
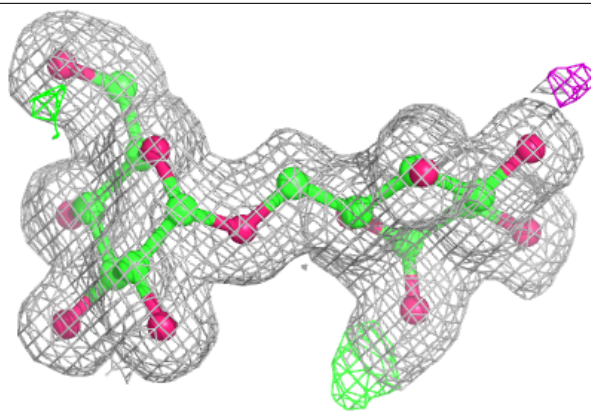
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GLC	L	3	11/12	0.74	0.24	44,45,47,47	0
3	GLC	L	4	11/12	0.74	0.28	43,49,51,51	0
3	GLC	M	1	11/12	0.77	0.17	27,36,39,41	0
3	GLC	K	4	11/12	0.79	0.26	36,40,41,42	0
3	GLC	K	3	11/12	0.79	0.34	41,44,46,48	0
4	GLC	G	2	11/12	0.80	0.26	29,31,32,34	11
4	GLC	G	3	11/12	0.80	0.28	34,34,35,35	11
3	GLC	L	1	11/12	0.82	0.16	37,41,43,43	0
3	GLC	L	2	11/12	0.83	0.17	40,42,43,44	0
4	GLC	G	1	12/12	0.83	0.23	29,31,31,32	12
3	GLC	K	2	11/12	0.83	0.31	42,43,43,43	0
3	GLC	K	1	11/12	0.83	0.23	33,37,38,40	0
3	GLC	F	3	11/12	0.84	0.15	21,27,29,29	0
2	GLC	I	1	12/12	0.85	0.14	27,35,37,38	0
3	GLC	D	3	11/12	0.85	0.16	36,40,44,46	0
3	GLC	M	4	11/12	0.85	0.11	34,36,38,40	0
5	GLC	H	3	11/12	0.86	0.17	29,30,31,31	0
3	GLC	M	3	11/12	0.88	0.10	24,25,28,31	0
3	GLC	D	1	11/12	0.88	0.11	32,34,36,39	0
2	GLC	I	2	11/12	0.90	0.09	22,26,28,28	0
3	GLC	D	2	11/12	0.91	0.16	41,42,45,46	0
3	GLC	F	2	11/12	0.91	0.22	28,30,32,32	0
5	GLC	H	4	11/12	0.91	0.14	22,26,27,27	0
2	GLC	E	1	12/12	0.92	0.12	22,30,33,34	0
5	GLC	H	2	11/12	0.92	0.09	22,22,25,27	0
3	GLC	F	1	11/12	0.93	0.08	18,19,21,25	0
5	GLC	H	1	11/12	0.94	0.08	17,18,20,20	0
3	GLC	D	4	11/12	0.94	0.10	31,32,34,34	0
2	GLC	J	1	12/12	0.94	0.09	18,21,21,22	0
2	GLC	E	2	11/12	0.94	0.07	20,21,21,22	0
2	GLC	C	2	11/12	0.96	0.07	15,17,17,18	0
2	GLC	C	1	12/12	0.97	0.06	17,20,22,23	0
3	GLC	F	4	11/12	0.97	0.07	15,16,18,18	0
3	GLC	M	2	11/12	0.97	0.07	19,20,23,23	0
2	GLC	J	2	11/12	0.97	0.05	16,16,17,17	0

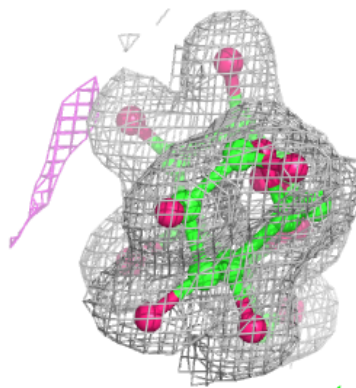
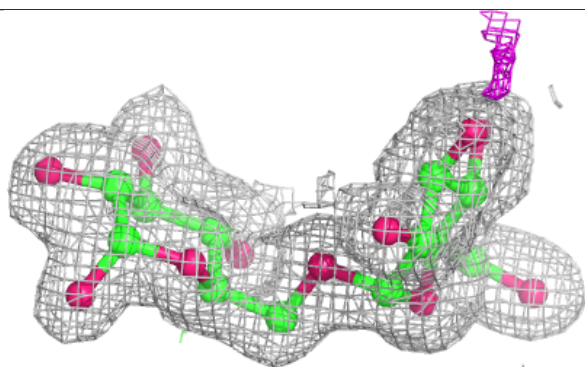
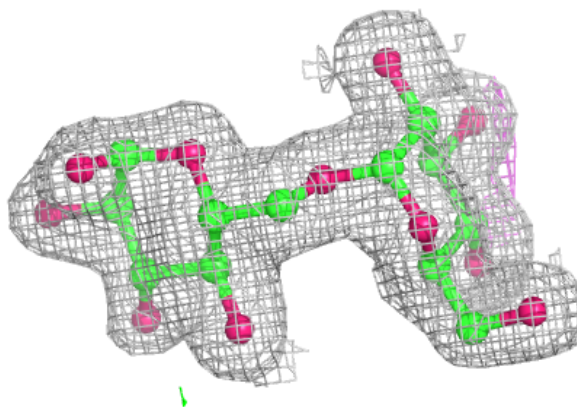
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

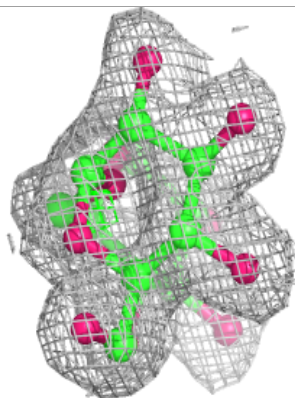
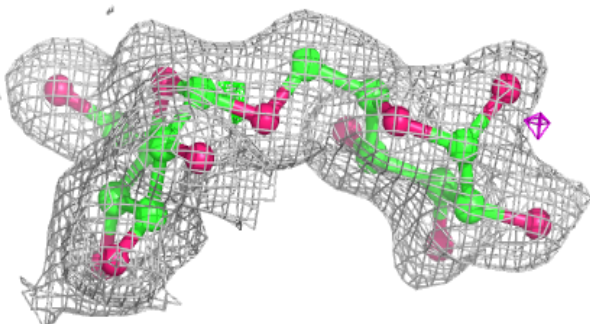
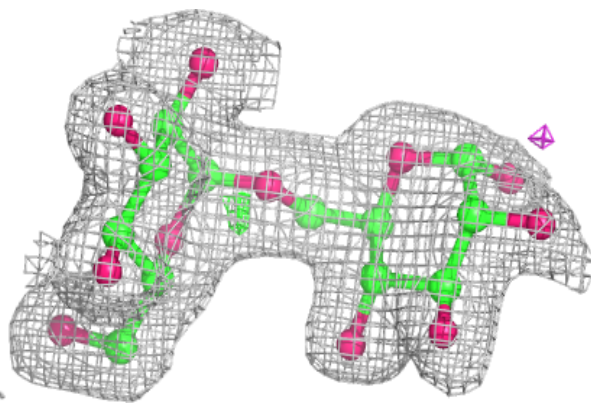
**Electron density around Chain E:**

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

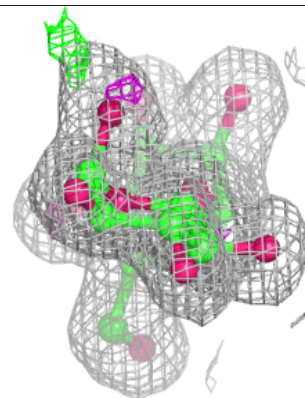
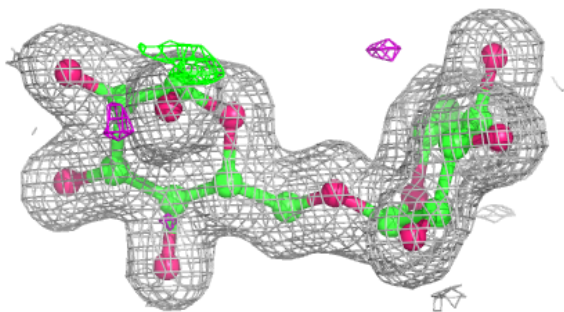
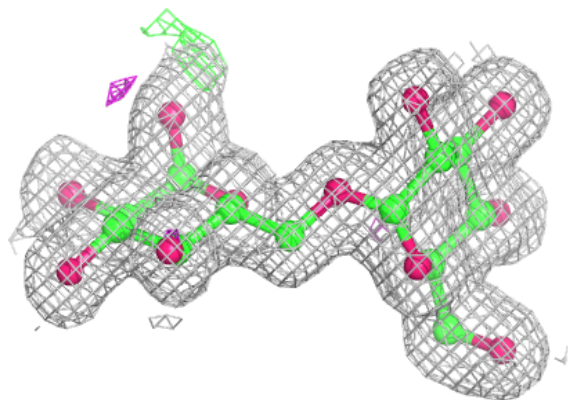


Electron density around Chain I:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

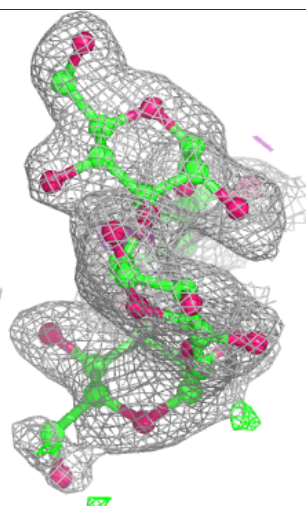
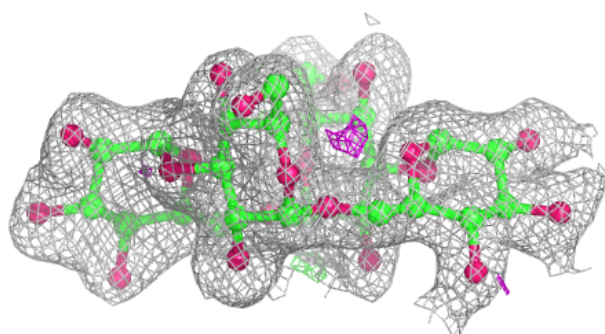
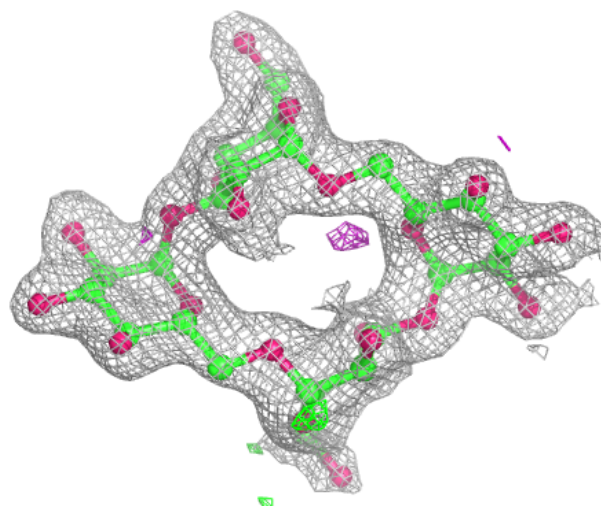
**Electron density around Chain J:**

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



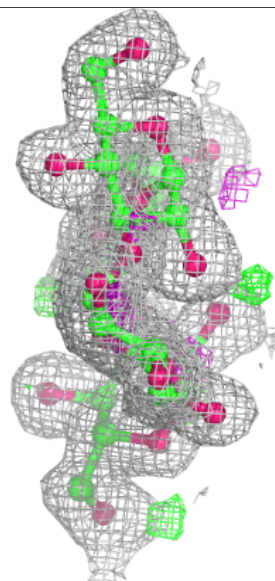
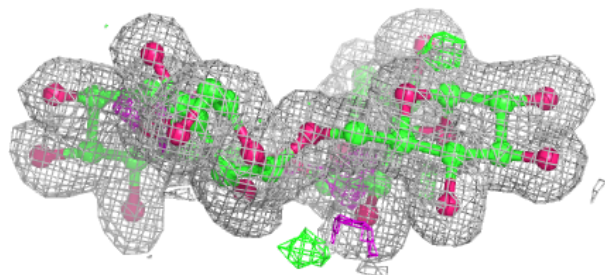
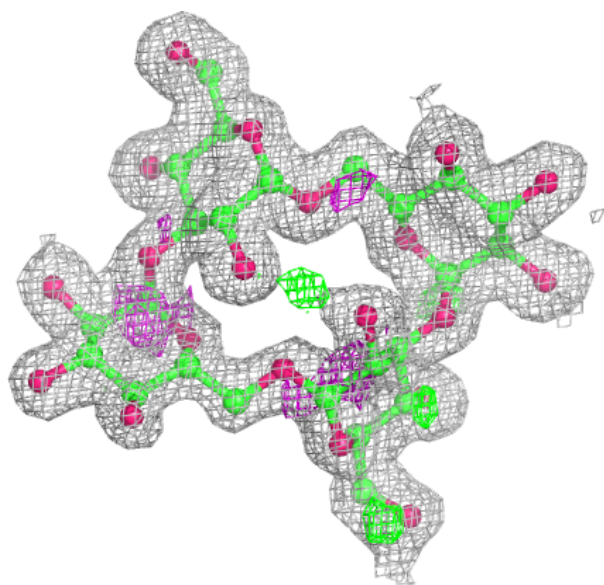
Electron density around Chain D:

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



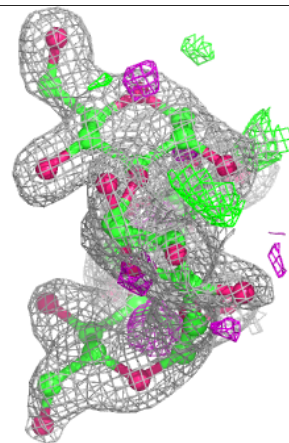
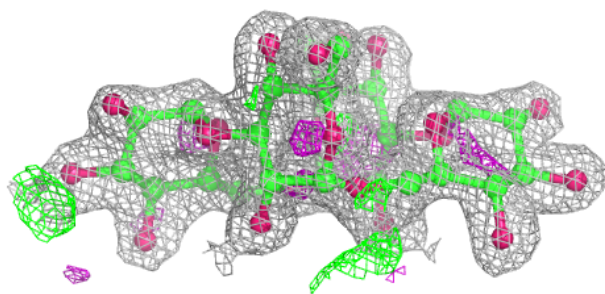
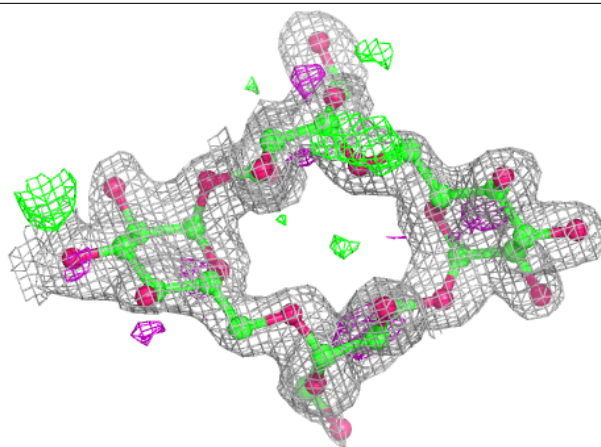
Electron density around Chain F:

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



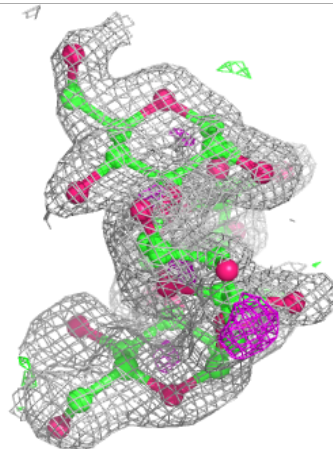
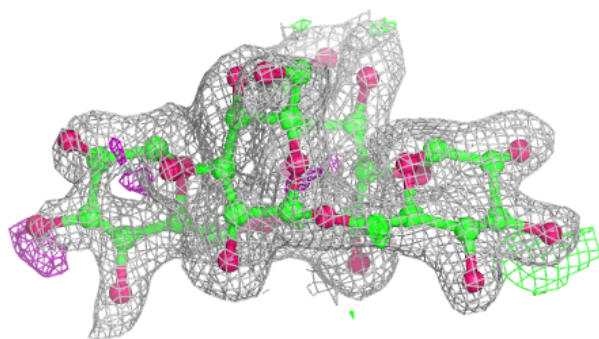
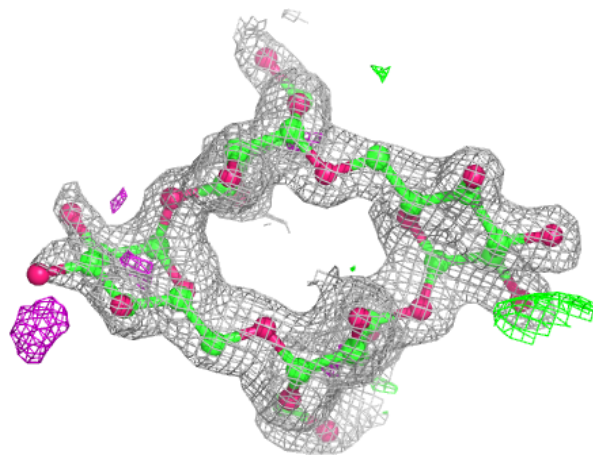
Electron density around Chain K:

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



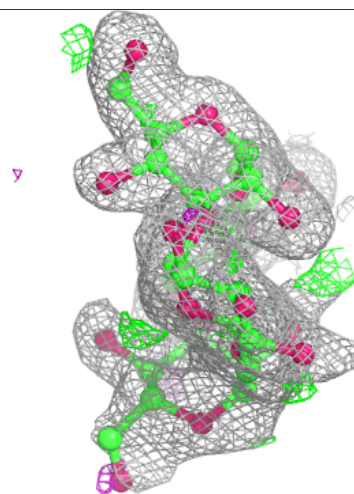
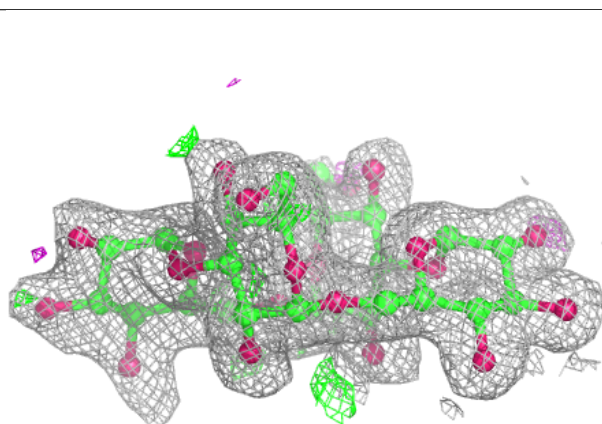
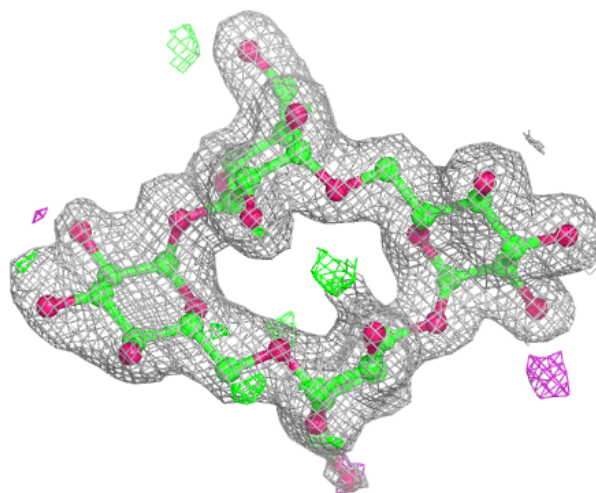
Electron density around Chain L:

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



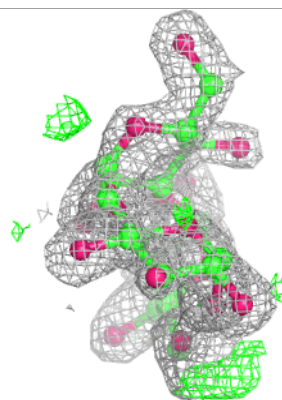
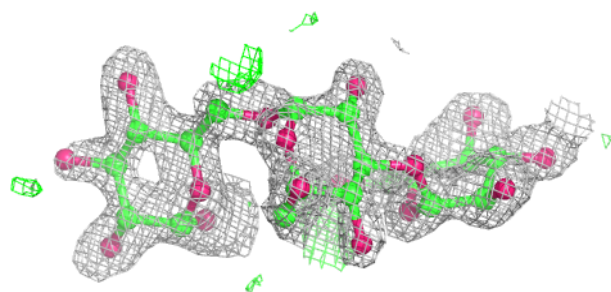
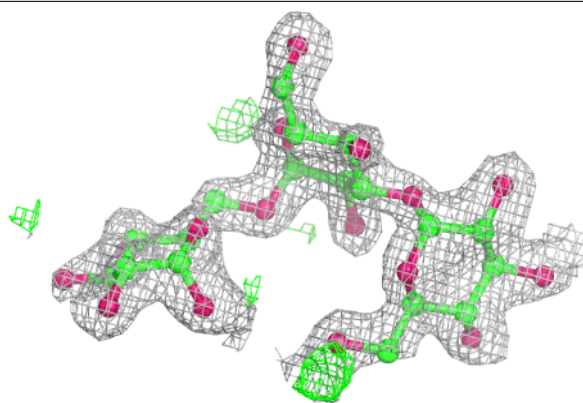
Electron density around Chain M:

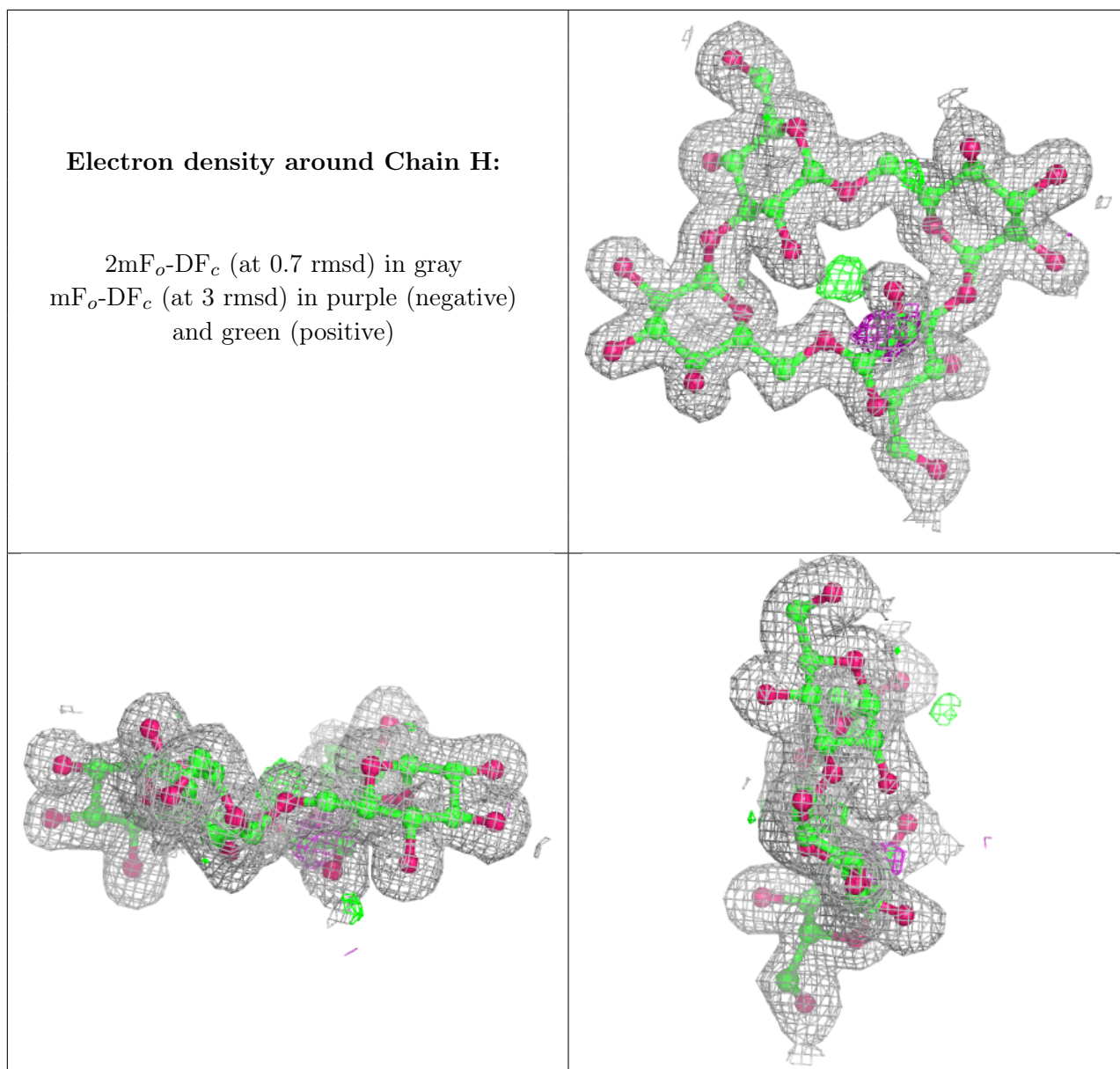
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain G:

$2mF_o-DF_c$ (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(Å ²)	Q<0.9
9	GLC	A	1124	12/12	0.75	0.20	29,31,33,35	12
9	GLC	B	1122	12/12	0.84	0.13	27,29,30,30	12
6	MG	A	1105	1/1	0.85	0.16	41,41,41,41	0
6	MG	A	1104	1/1	0.90	0.16	33,33,33,33	0
8	CL	B	1108	1/1	0.94	0.15	34,34,34,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
8	CL	A	1108	1/1	0.96	0.06	40,40,40,40	0
6	MG	B	1102	1/1	0.96	0.07	17,17,17,17	0
8	CL	B	1105	1/1	0.97	0.04	26,26,26,26	0
6	MG	A	1102	1/1	0.97	0.13	25,25,25,25	0
6	MG	A	1101	1/1	0.98	0.04	14,14,14,14	0
8	CL	B	1107	1/1	0.98	0.04	33,33,33,33	0
6	MG	B	1103	1/1	0.98	0.09	32,32,32,32	0
7	CA	A	1106	1/1	0.98	0.03	14,14,14,14	0
6	MG	A	1103	1/1	0.98	0.05	14,14,14,14	0
8	CL	A	1107	1/1	0.99	0.08	35,35,35,35	0
6	MG	B	1101	1/1	0.99	0.04	15,15,15,15	0
8	CL	B	1109	1/1	0.99	0.03	24,24,24,24	0
7	CA	B	1104	1/1	0.99	0.04	11,11,11,11	0
8	CL	B	1106	1/1	0.99	0.06	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.