



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 13, 2024 – 09:48 pm GMT

PDB ID : 6TIG  
Title : Structure of the N terminal domain of Bc2L-C lectin (1-131) in complex with Globo H (H-type 3) antigen  
Authors : Varrot, A.; Bermeo, R.  
Deposited on : 2019-11-22  
Resolution : 1.90 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](mailto: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 : 1.8.4, 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

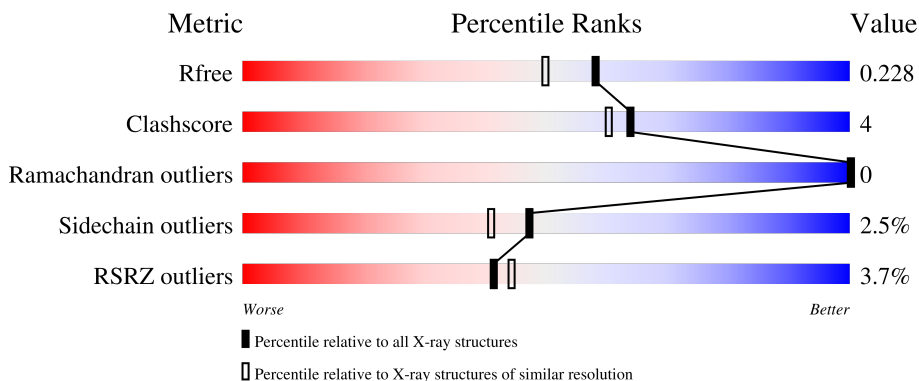
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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  $\geq 3$ , 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	AAA	134	4% 90% 10%
1	BBB	134	3% 87% 13%
1	CCC	134	4% 87% 13%
2	A	4	100%
3	B	5	100%

Continued on next page...

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
3	C	5	 100%

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 3483 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.

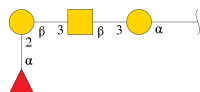
- Molecule 1 is a protein called Lectin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	AAA	134	Total 994	C 639	N 162	O 189	S 4	0	2	0
1	BBB	134	Total 1027	C 662	N 166	O 195	S 4	0	5	0
1	CCC	134	Total 1002	C 645	N 162	O 191	S 4	0	3	0

There are 6 discrepancies between the modelled and reference sequences:

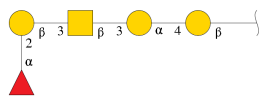
Chain	Residue	Modelled	Actual	Comment	Reference
AAA	-2	GLY	-	expression tag	UNP B4EH86
AAA	-1	HIS	-	expression tag	UNP B4EH86
BBB	-2	GLY	-	expression tag	UNP B4EH86
BBB	-1	HIS	-	expression tag	UNP B4EH86
CCC	-2	GLY	-	expression tag	UNP B4EH86
CCC	-1	HIS	-	expression tag	UNP B4EH86

- Molecule 2 is an oligosaccharide called alpha-L-fucopyranose-(1-2)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-galactopyranose-(1-3)-alpha-D-galactopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	A	4	Total 47	C 26	N 1	O 20	0	0	0

- Molecule 3 is an oligosaccharide called alpha-L-fucopyranose-(1-2)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-galactopyranose-(1-3)-alpha-D-galactopyranose-(1-4)-beta-D-galactopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	B	5	58	32	1	25	0	0	0
3	C	5	58	32	1	25	0	0	0

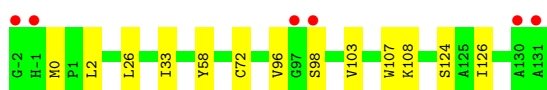
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	AAA	104	109	109	0	5
4	BBB	103	106	106	0	3
4	CCC	80	82	82	0	2

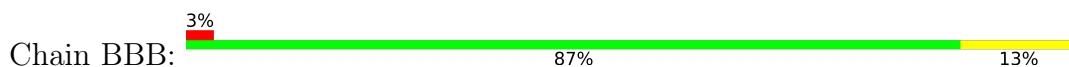
### 3 Residue-property plots [i](#)

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.

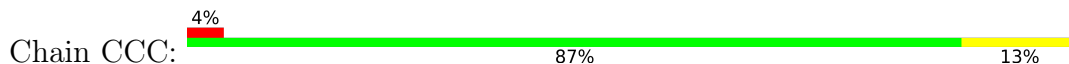
- Molecule 1: Lectin



- Molecule 1: Lectin



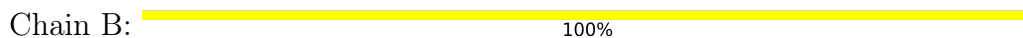
- Molecule 1: Lectin



- Molecule 2: alpha-L-fucopyranose-(1-2)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-galactopyranose-(1-3)-alpha-D-galactopyranose



- Molecule 3: alpha-L-fucopyranose-(1-2)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-galactopyranose-(1-3)-alpha-D-galactopyranose-(1-4)-beta-D-galactopyranose



- Molecule 3: alpha-L-fucopyranose-(1-2)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-galactopyranose-(1-3)-alpha-D-galactopyranose-(1-4)-beta-D-galactopyranose

Chain C:

100%

GAL1  
GAL2  
NGA3  
GAL4  
FUC5

## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.37Å 42.88Å 102.65Å 90.00° 96.01° 90.00°	Depositor
Resolution (Å)	37.10 – 1.90 37.10 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.5 (37.10-1.90) 97.6 (37.10-1.90)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.32 (at 1.89Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.165 , 0.227 0.176 , 0.228	Depositor DCC
$R_{free}$ test set	1552 reflections (6.19%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	21.4	Xtrriage
Anisotropy	0.563	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 50.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3483	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.43% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: NGA, GAL, GLA, FUC

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	AAA	0.85	0/1018	0.95	0/1393
1	BBB	0.83	0/1055	0.95	0/1441
1	CCC	0.89	1/1029 (0.1%)	0.99	1/1408 (0.1%)
All	All	0.86	1/3102 (0.0%)	0.96	1/4242 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	CCC	81	GLU	CD-OE1	-6.12	1.19	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CCC	111	ARG	NE-CZ-NH1	-5.54	117.53	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AAA	994	0	988	12	0
1	BBB	1027	0	1027	12	0
1	CCC	1002	0	1001	8	1

*Continued on next page...*

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	47	0	42	0	0
3	B	58	0	51	0	0
3	C	58	0	51	0	0
4	AAA	109	0	0	0	0
4	BBB	106	0	0	1	0
4	CCC	82	0	0	0	0
All	All	3483	0	3160	27	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 4.

All (27) 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:BBB:70[B]:ASP:OD2	4:BBB:301:HOH:O	1.74	1.04
1:AAA:26:LEU:HB2	1:AAA:103[B]:VAL:HG22	1.75	0.68
1:CCC:2:LEU:HD22	1:CCC:126:ILE:HG12	1.80	0.63
1:CCC:26:LEU:HB2	1:CCC:103[B]:VAL:HG22	1.82	0.62
1:BBB:2:LEU:HD22	1:BBB:126:ILE:HG12	1.82	0.62
1:BBB:26:LEU:HB2	1:BBB:103[A]:VAL:HG22	1.81	0.62
1:CCC:33:ILE:HG22	1:CCC:96:VAL:HG11	1.84	0.60
1:AAA:0:MET:HG2	1:BBB:0:MET:HE1	1.85	0.59
1:BBB:33:ILE:HG22	1:BBB:96:VAL:HG11	1.86	0.58
1:AAA:33:ILE:HG22	1:AAA:96:VAL:HG11	1.87	0.54
1:AAA:0:MET:CE	1:CCC:0:MET:HG2	2.39	0.52
1:CCC:37:LYS:HE2	1:CCC:95:ASP:OD1	2.11	0.51
1:AAA:26:LEU:HB2	1:AAA:103[B]:VAL:CG2	2.42	0.49
1:BBB:54:PHE:CE1	1:CCC:82:SER:HB3	2.46	0.49
1:AAA:0:MET:HE1	1:CCC:0:MET:HG2	1.96	0.46
1:AAA:26:LEU:HD12	1:AAA:103[B]:VAL:CG2	2.46	0.46
1:BBB:53:ASN:OD1	1:BBB:78[A]:LYS:NZ	2.45	0.45
1:AAA:2:LEU:HD22	1:AAA:126:ILE:HG12	1.99	0.44
1:AAA:126:ILE:HG21	1:BBB:126:ILE:HD13	1.99	0.44
1:BBB:58[A]:TYR:CD1	1:BBB:72:CYS:HB3	2.54	0.43
1:BBB:58[A]:TYR:HB2	1:BBB:108:LYS:HG2	1.99	0.43
1:BBB:26:LEU:HD12	1:BBB:103[A]:VAL:CG2	2.49	0.43
1:AAA:58:TYR:CD1	1:AAA:72:CYS:HB3	2.55	0.42
1:CCC:58:TYR:CD1	1:CCC:72:CYS:HB3	2.54	0.42
1:AAA:58:TYR:HB2	1:AAA:108:LYS:HG2	2.02	0.41
1:AAA:26:LEU:HD12	1:AAA:103[B]:VAL:HG21	2.03	0.41
1:BBB:57:ILE:O	1:BBB:72:CYS:HB2	2.21	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)
1:CCC:32:GLY:O	1:CCC:32:GLY:O[2_556]	2.18	0.02

### 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	AAA	134/134 (100%)	129 (96%)	5 (4%)	0	100	100
1	BBB	137/134 (102%)	133 (97%)	4 (3%)	0	100	100
1	CCC	135/134 (101%)	131 (97%)	4 (3%)	0	100	100
All	All	406/402 (101%)	393 (97%)	13 (3%)	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	AAA	106/106 (100%)	103 (97%)	3 (3%)	43	36
1	BBB	111/106 (105%)	109 (98%)	2 (2%)	59	55
1	CCC	108/106 (102%)	105 (97%)	3 (3%)	43	36
All	All	325/318 (102%)	317 (98%)	8 (2%)	47	41

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	98	SER
1	AAA	107	TRP
1	AAA	124	SER
1	BBB	98	SER
1	BBB	107	TRP
1	CCC	98	SER
1	CCC	107	TRP
1	CCC	124	SER

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

### 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](#)

14 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	GLA	A	1	2	12,12,12	1.05	1 (8%)	17,17,17	1.59	5 (29%)
2	NGA	A	2	2	14,14,15	1.01	1 (7%)	17,19,21	2.29	8 (47%)
2	GAL	A	3	2	11,11,12	1.27	2 (18%)	15,15,17	1.66	6 (40%)
2	FUC	A	4	2	10,10,11	1.71	3 (30%)	14,14,16	1.13	1 (7%)
3	GAL	B	1	3	12,12,12	0.84	0	17,17,17	1.36	3 (17%)
3	GLA	B	2	3	11,11,12	1.19	1 (9%)	15,15,17	3.54	10 (66%)
3	NGA	B	3	3	14,14,15	0.85	0	17,19,21	1.50	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	GAL	B	4	3	11,11,12	1.09	1 (9%)	15,15,17	2.24	6 (40%)
3	FUC	B	5	3	10,10,11	1.44	2 (20%)	14,14,16	1.57	4 (28%)
3	GAL	C	1	3	12,12,12	1.10	1 (8%)	17,17,17	1.86	4 (23%)
3	GLA	C	2	3	11,11,12	1.20	1 (9%)	15,15,17	2.09	5 (33%)
3	NGA	C	3	3	14,14,15	1.09	1 (7%)	17,19,21	2.17	5 (29%)
3	GAL	C	4	3	11,11,12	1.23	1 (9%)	15,15,17	1.58	3 (20%)
3	FUC	C	5	3	10,10,11	1.32	1 (10%)	14,14,16	1.36	2 (14%)

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	GLA	A	1	2	-	0/2/22/22	0/1/1/1
2	NGA	A	2	2	-	1/6/23/26	0/1/1/1
2	GAL	A	3	2	-	2/2/19/22	0/1/1/1
2	FUC	A	4	2	-	-	0/1/1/1
3	GAL	B	1	3	-	1/2/22/22	0/1/1/1
3	GLA	B	2	3	-	0/2/19/22	0/1/1/1
3	NGA	B	3	3	-	2/6/23/26	0/1/1/1
3	GAL	B	4	3	-	1/2/19/22	0/1/1/1
3	FUC	B	5	3	-	-	0/1/1/1
3	GAL	C	1	3	-	1/2/22/22	0/1/1/1
3	GLA	C	2	3	-	2/2/19/22	0/1/1/1
3	NGA	C	3	3	-	0/6/23/26	0/1/1/1
3	GAL	C	4	3	-	1/2/19/22	0/1/1/1
3	FUC	C	5	3	-	-	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	4	FUC	C2-C3	-3.89	1.46	1.52
3	C	5	FUC	O5-C5	-3.48	1.36	1.43
3	C	2	GLA	C2-C3	3.22	1.57	1.52
3	B	2	GLA	C2-C3	2.99	1.56	1.52
3	B	5	FUC	O5-C5	-2.82	1.37	1.43
3	B	4	GAL	O2-C2	2.60	1.48	1.43
3	C	1	GAL	C1-C2	2.51	1.58	1.52
3	C	4	GAL	O5-C5	-2.50	1.38	1.43

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	3	GAL	C4-C5	-2.27	1.48	1.53
3	B	5	FUC	C4-C3	2.25	1.58	1.52
3	C	3	NGA	C4-C5	2.25	1.57	1.53
2	A	4	FUC	O2-C2	-2.24	1.38	1.43
2	A	3	GAL	O5-C1	-2.24	1.40	1.43
2	A	4	FUC	C4-C3	2.23	1.58	1.52
2	A	2	NGA	C1-C2	2.08	1.55	1.52
2	A	1	GLA	C1-C2	2.07	1.57	1.52

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	2	GLA	O4-C4-C3	-5.44	97.78	110.35
3	B	2	GLA	O2-C2-C3	5.44	121.03	110.14
3	C	2	GLA	O3-C3-C4	-5.21	98.31	110.35
3	B	2	GLA	C6-C5-C4	4.88	124.44	113.00
3	B	2	GLA	O4-C4-C5	4.87	121.38	109.30
3	B	2	GLA	C3-C4-C5	-4.64	101.95	110.24
3	C	3	NGA	C1-O5-C5	-4.61	105.94	112.19
3	B	4	GAL	C1-C2-C3	-4.59	104.03	109.67
3	B	2	GLA	O3-C3-C2	4.33	118.28	109.99
3	B	4	GAL	O5-C1-C2	4.05	117.02	110.77
3	C	3	NGA	O4-C4-C5	4.04	119.32	109.30
2	A	2	NGA	C2-N2-C7	-3.95	117.27	122.90
3	C	1	GAL	C4-C3-C2	-3.76	104.25	110.82
3	B	2	GLA	C2-C3-C4	-3.74	104.42	110.89
2	A	2	NGA	C1-O5-C5	-3.73	107.13	112.19
2	A	2	NGA	O5-C5-C6	3.46	112.63	107.20
3	C	3	NGA	C1-C2-N2	-3.22	104.98	110.49
3	C	1	GAL	O2-C2-C1	3.14	116.44	109.16
2	A	2	NGA	C1-C2-N2	-3.10	105.20	110.49
3	C	5	FUC	C2-C3-C4	-3.06	105.61	110.89
3	B	3	NGA	O3-C3-C4	-3.04	103.31	110.35
2	A	1	GLA	O2-C2-C3	-3.01	103.40	110.35
3	B	2	GLA	O3-C3-C4	-2.99	103.44	110.35
3	B	4	GAL	C2-C3-C4	2.98	116.06	110.89
2	A	3	GAL	O4-C4-C3	2.90	117.05	110.35
3	C	2	GLA	O5-C5-C6	2.87	111.70	107.20
3	B	4	GAL	O4-C4-C5	2.81	116.28	109.30
3	B	5	FUC	C2-C3-C4	-2.70	106.23	110.89
3	C	2	GLA	O3-C3-C2	2.68	115.12	109.99
3	B	4	GAL	C3-C4-C5	-2.66	105.50	110.24

*Continued on next page...*

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	3	GAL	O3-C3-C4	-2.64	104.25	110.35
2	A	2	NGA	O7-C7-N2	-2.63	117.11	121.95
3	B	3	NGA	C1-O5-C5	-2.59	108.69	112.19
2	A	1	GLA	O1-C1-C2	2.58	116.30	109.03
3	C	3	NGA	O5-C1-C2	2.54	115.30	111.29
3	B	1	GAL	O5-C5-C6	2.54	112.74	106.44
3	C	4	GAL	O4-C4-C3	2.49	116.11	110.35
2	A	1	GLA	O1-C1-O5	-2.48	102.93	110.38
3	B	1	GAL	O3-C3-C2	2.48	116.08	110.35
2	A	2	NGA	O7-C7-C8	2.48	126.66	122.06
3	B	3	NGA	O5-C5-C6	-2.47	103.33	107.20
2	A	4	FUC	C2-C3-C4	-2.43	106.69	110.89
3	C	1	GAL	O4-C4-C3	-2.42	104.75	110.35
2	A	3	GAL	O2-C2-C3	-2.40	105.34	110.14
3	B	2	GLA	C1-O5-C5	-2.39	108.95	112.19
3	B	5	FUC	O4-C4-C5	2.36	114.90	109.67
3	B	3	NGA	C2-N2-C7	-2.36	119.54	122.90
3	B	4	GAL	C1-O5-C5	-2.34	109.02	112.19
3	B	1	GAL	O4-C4-C5	2.33	115.08	109.30
2	A	2	NGA	O4-C4-C5	-2.30	103.59	109.30
3	C	2	GLA	O2-C2-C3	2.30	114.74	110.14
3	C	5	FUC	O5-C5-C6	-2.29	102.39	107.33
2	A	2	NGA	O5-C5-C4	-2.26	105.32	110.83
3	C	4	GAL	O5-C1-C2	2.24	114.24	110.77
2	A	3	GAL	O3-C3-C2	2.21	114.22	109.99
3	C	2	GLA	O5-C1-C2	-2.19	107.39	110.77
3	B	2	GLA	O6-C6-C5	-2.19	103.78	111.29
3	C	1	GAL	O1-C1-O5	-2.17	103.86	110.38
2	A	3	GAL	C1-O5-C5	-2.17	109.25	112.19
2	A	1	GLA	O2-C2-C1	2.15	114.15	109.16
3	B	5	FUC	O5-C5-C4	-2.11	105.73	109.52
3	C	3	NGA	O5-C5-C6	2.11	110.51	107.20
2	A	3	GAL	C1-C2-C3	-2.10	107.09	109.67
3	C	4	GAL	O2-C2-C1	-2.03	104.99	109.15
2	A	1	GLA	O5-C5-C4	2.02	113.36	109.69
3	B	5	FUC	C3-C4-C5	2.01	112.91	109.77

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	3	NGA	O5-C5-C6-O6

Continued on next page...

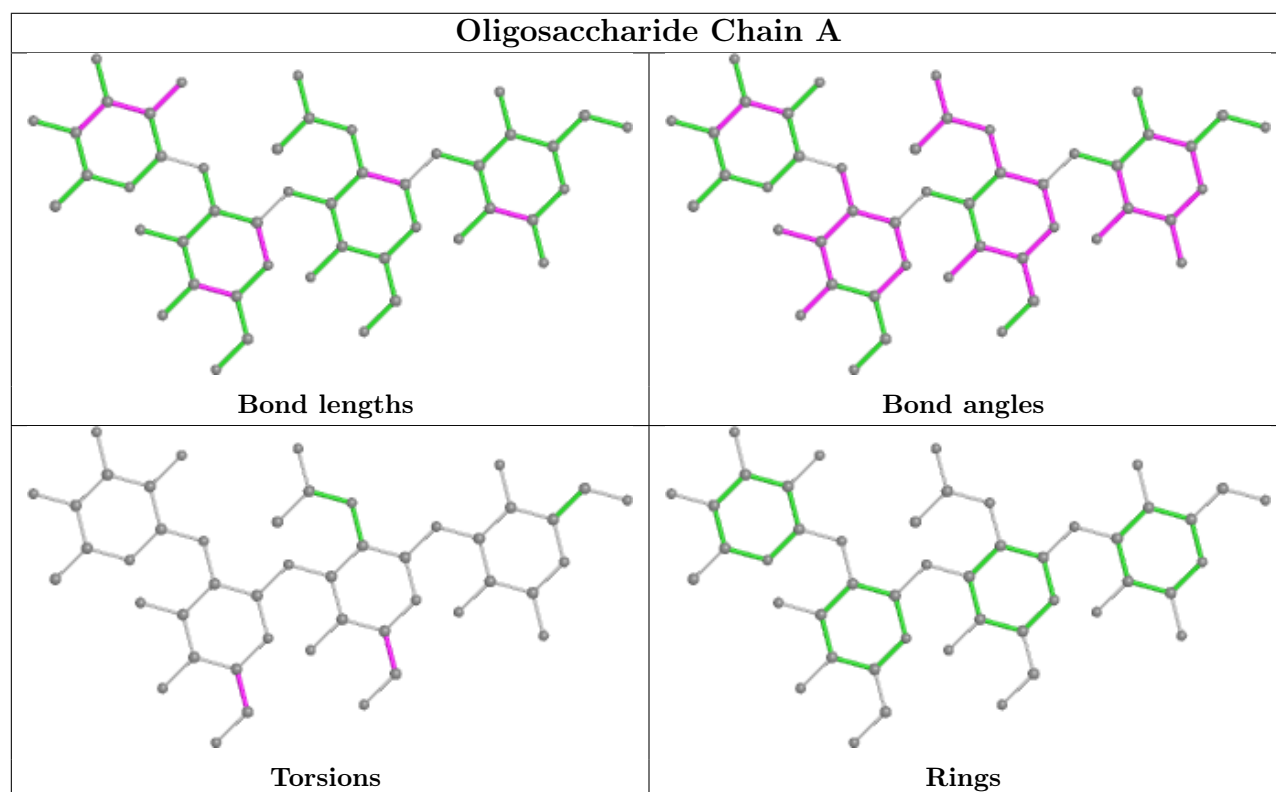
Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	C	2	GLA	C4-C5-C6-O6
3	C	2	GLA	O5-C5-C6-O6
3	B	3	NGA	C4-C5-C6-O6
3	B	4	GAL	C4-C5-C6-O6
3	B	1	GAL	O5-C5-C6-O6
3	C	1	GAL	O5-C5-C6-O6
2	A	3	GAL	O5-C5-C6-O6
3	C	4	GAL	C4-C5-C6-O6
2	A	3	GAL	C4-C5-C6-O6
2	A	2	NGA	C4-C5-C6-O6

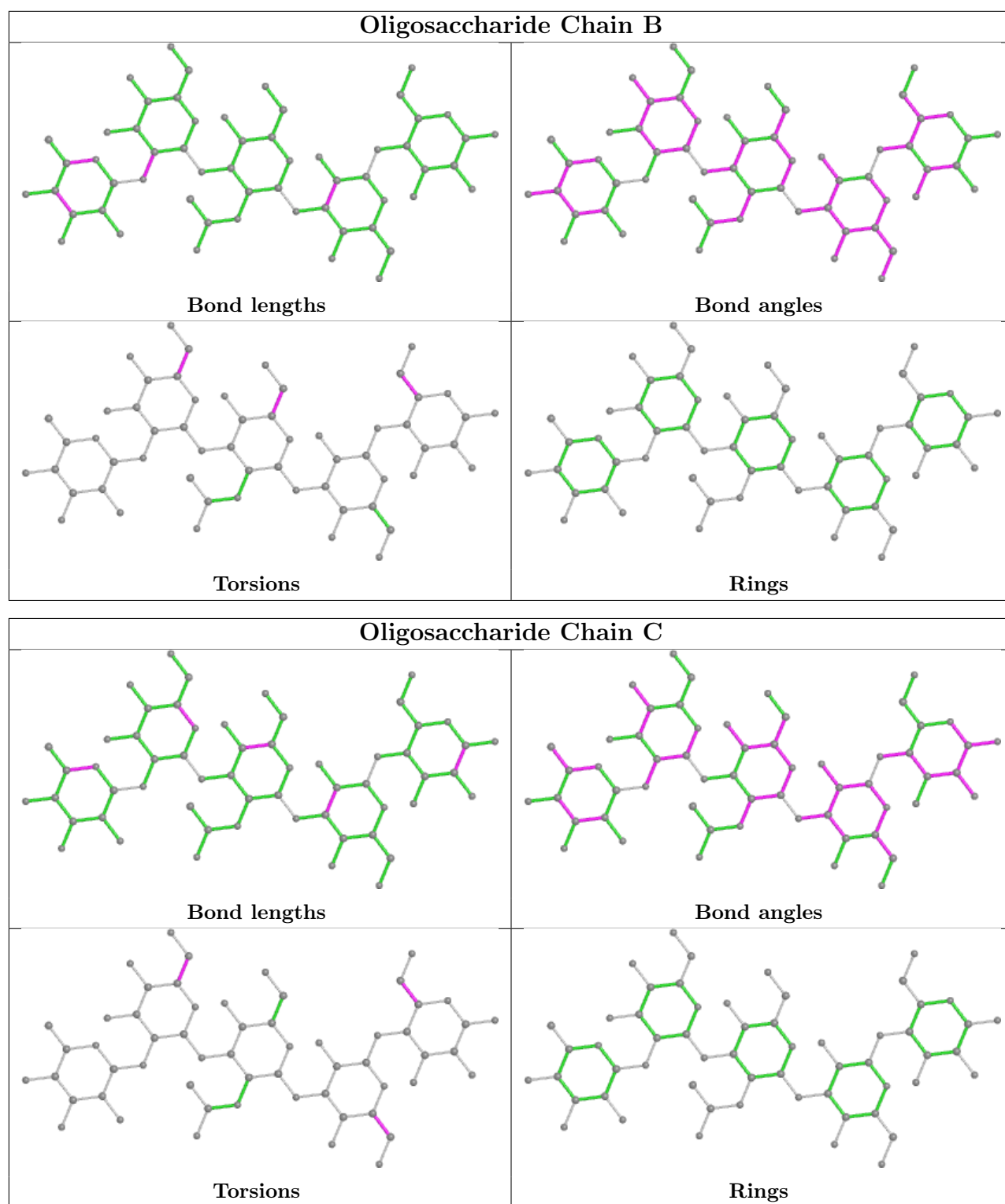
There are no ring outliers.

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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	AAA	134/134 (100%)	0.03	6 (4%) 33 36	15, 22, 49, 63	0
1	BBB	134/134 (100%)	0.06	4 (2%) 50 53	15, 22, 50, 64	1 (0%)
1	CCC	134/134 (100%)	0.12	5 (3%) 41 44	15, 22, 50, 63	1 (0%)
All	All	402/402 (100%)	0.07	15 (3%) 41 44	15, 22, 50, 64	2 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	131	ALA	5.2
1	CCC	131	ALA	3.6
1	CCC	97	GLY	3.3
1	CCC	29	ALA	3.2
1	AAA	130	ALA	3.0
1	AAA	-1	HIS	2.8
1	BBB	-2	GLY	2.7
1	AAA	98	SER	2.6
1	CCC	98	SER	2.5
1	BBB	98	SER	2.5
1	AAA	97	GLY	2.5
1	BBB	131	ALA	2.4
1	BBB	97	GLY	2.3
1	AAA	-2	GLY	2.2
1	CCC	127	TRP	2.2

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates

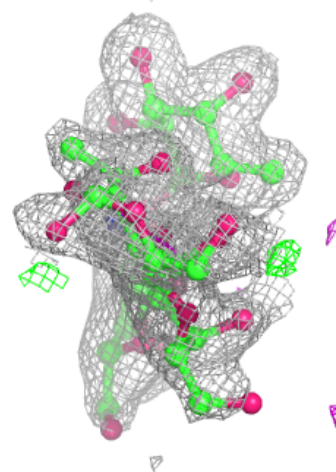
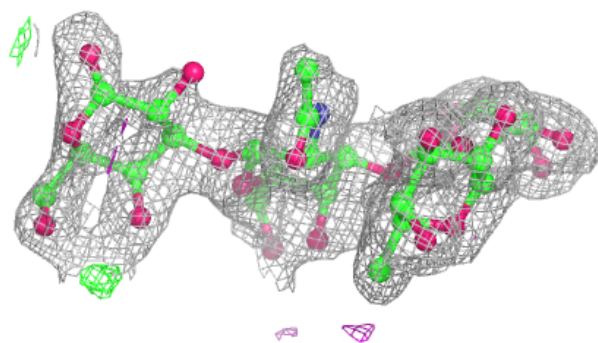
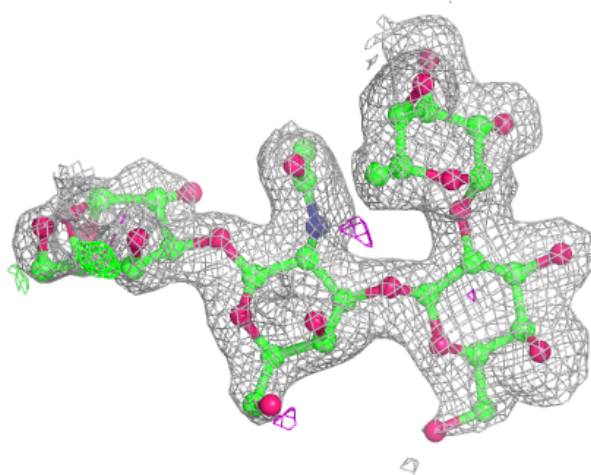
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	GLA	B	2	11/12	0.74	0.17	49,59,66,67	0
3	GAL	B	1	12/12	0.78	0.20	50,53,65,66	0
2	GLA	A	1	12/12	0.78	0.29	36,46,50,50	0
3	GAL	C	1	12/12	0.83	0.16	41,48,53,57	0
3	GLA	C	2	11/12	0.90	0.29	39,49,58,61	0
3	NGA	C	3	14/15	0.90	0.19	28,32,47,52	0
3	NGA	B	3	14/15	0.92	0.15	29,34,43,54	0
3	GAL	C	4	11/12	0.92	0.14	24,32,44,45	0
2	GAL	A	3	11/12	0.93	0.11	21,29,41,47	0
2	NGA	A	2	14/15	0.94	0.18	25,30,46,50	0
3	GAL	B	4	11/12	0.94	0.12	23,31,43,44	0
2	FUC	A	4	10/11	0.97	0.07	17,19,21,22	0
3	FUC	B	5	10/11	0.97	0.08	17,20,22,22	0
3	FUC	C	5	10/11	0.98	0.06	17,19,22,22	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.

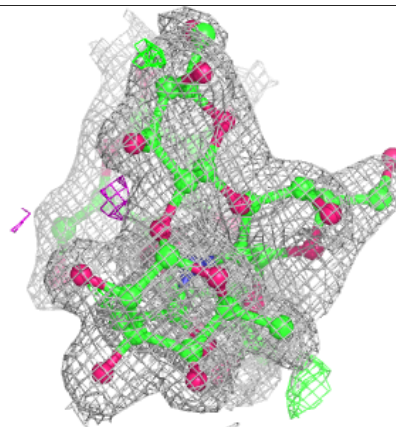
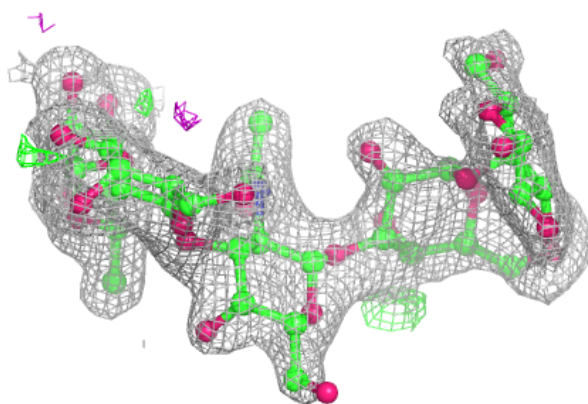
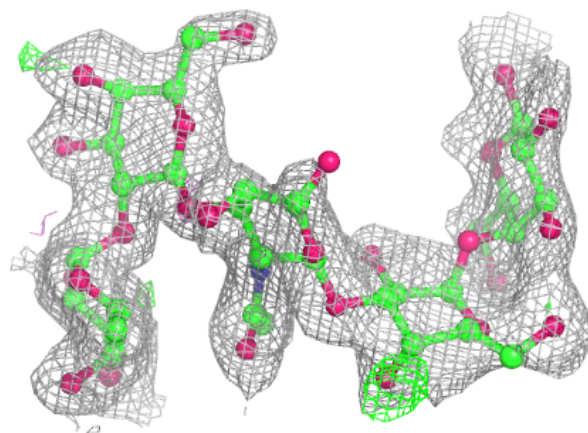
**Electron density around Chain A:**

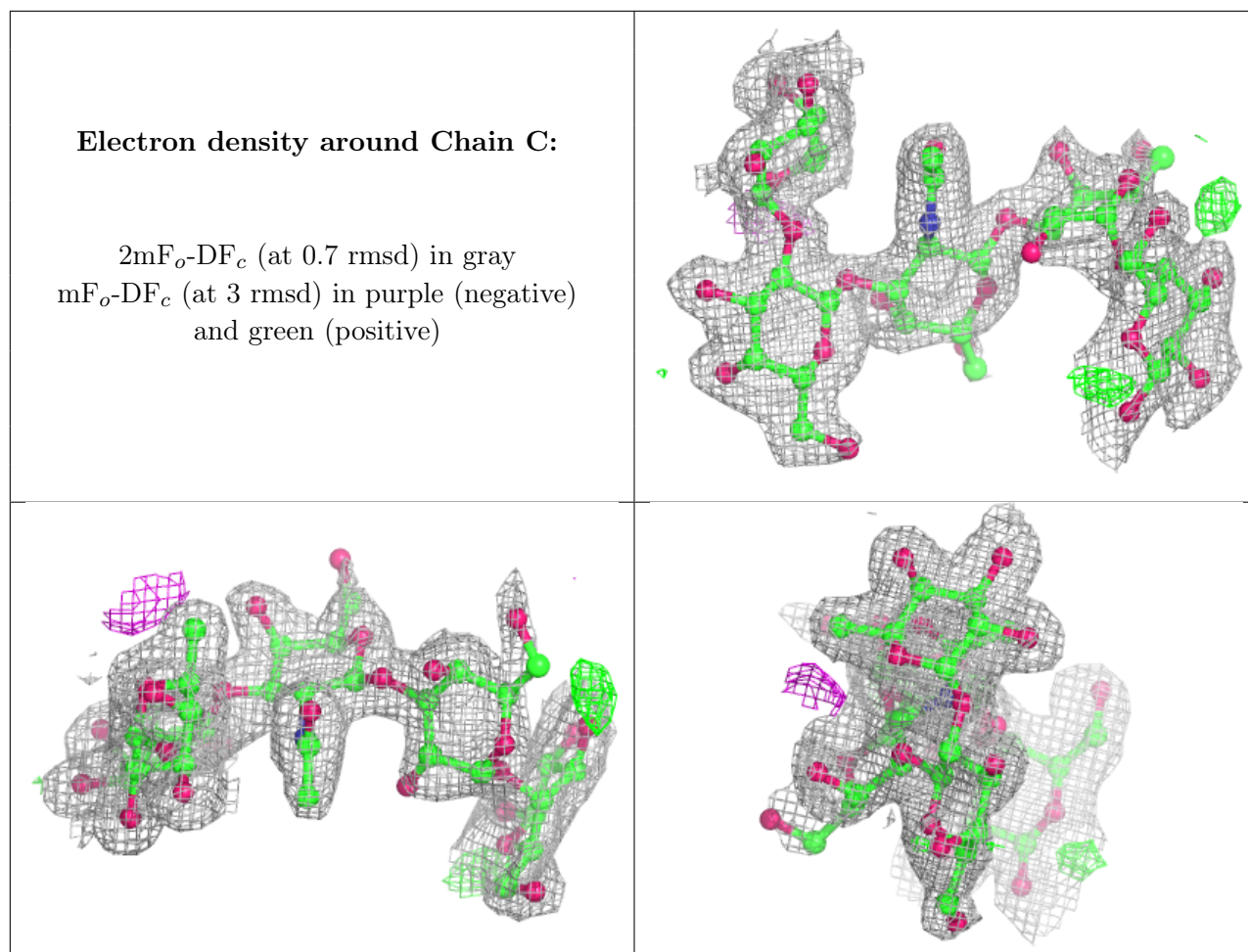
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around Chain B:**

$2mF_o-DF_c$  (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.