



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 23, 2024 – 11:51 PM EDT

PDB ID : 2OVW  
Title : ENDOGLUCANASE I COMPLEXED WITH CELLOBIOSE  
Authors : Sulzenbacher, G.; Davies, G.J.; Schulein, M.  
Deposited on : 1997-04-04  
Resolution : 2.30 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](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>.

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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 : 2.36.1  
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

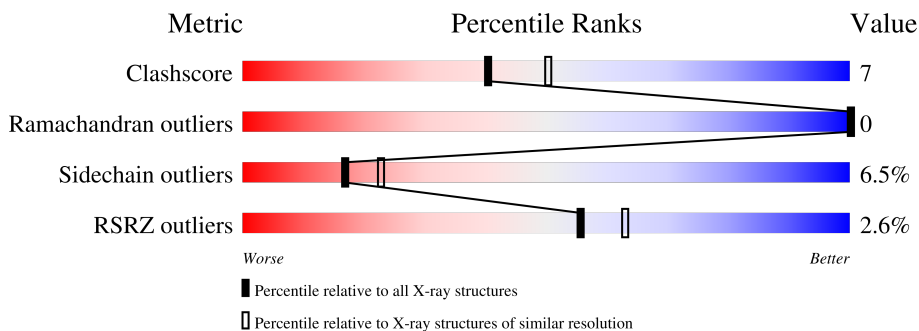
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	A	411	 2% 75% 19% . .
1	B	411	 4% 75% 18% . .
1	C	411	 2% 74% 19% . .
1	D	411	 % 72% 21% . .
2	E	2	 50% 50%
2	F	2	 50% 50%

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Mol	Chain	Length	Quality of chain
2	G	2	 50% 50%
2	H	2	 50% 50%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 13393 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 ENDOGLUCANASE I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	398	3022	1870	534	589	29	0	0	0
1	B	398	3022	1870	534	589	29	0	0	0
1	C	398	3022	1870	534	589	29	0	0	0
1	D	398	3022	1870	534	589	29	0	0	0

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	E	2	23	12	11	0	0	0
2	F	2	23	12	11	0	0	0
2	G	2	23	12	11	0	0	0
2	H	2	23	12	11	0	0	0

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	14	8	1	5	0	0
3	A	1	14	8	1	5	0	0
3	B	1	14	8	1	5	0	0
3	B	1	14	8	1	5	0	0
3	C	1	14	8	1	5	0	0
3	C	1	14	8	1	5	0	0
3	D	1	14	8	1	5	0	0
3	D	1	14	8	1	5	0	0

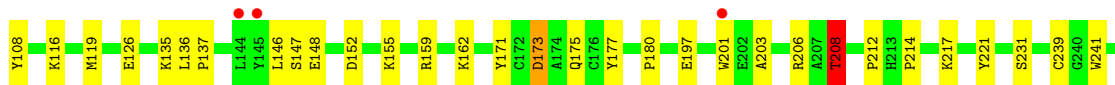
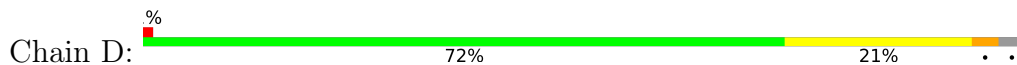
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	277	Total 277	O 277	0	0
4	B	268	Total 268	O 268	0	0
4	C	268	Total 268	O 268	0	0
4	D	288	Total 288	O 288	0	0





• Molecule 1: ENDOGLUCANASE I



ALA	PRO
ALA	TYR
PRO	GLY
PRO	HIS
ARG	LEU

• Molecule 2: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



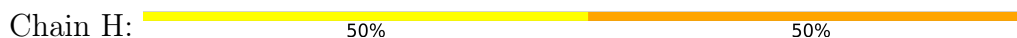
• Molecule 2: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



• Molecule 2: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



• Molecule 2: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.93Å 78.31Å 142.47Å 90.00° 96.98° 90.00°	Depositor
Resolution (Å)	15.00 – 2.30 29.84 – 2.30	Depositor EDS
% Data completeness (in resolution range)	86.4 (15.00-2.30) 86.5 (29.84-2.30)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.29 (at 2.31Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.209 , (Not available) 0.195 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.4	Xtrriage
Anisotropy	0.487	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 61.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13393	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.25% 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: BGC, PCA, NAG

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	0.68	0/3076	1.58	41/4155 (1.0%)
1	B	0.68	0/3076	1.59	40/4155 (1.0%)
1	C	0.70	0/3076	1.57	34/4155 (0.8%)
1	D	0.69	0/3076	1.59	44/4155 (1.1%)
All	All	0.69	0/12304	1.58	159/16620 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	4

There are no bond length outliers.

The worst 5 of 159 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	325	ARG	NE-CZ-NH1	18.48	129.54	120.30
1	A	325	ARG	NE-CZ-NH1	16.49	128.55	120.30
1	C	264	ARG	CD-NE-CZ	15.53	145.34	123.60
1	C	325	ARG	CD-NE-CZ	15.48	145.27	123.60
1	C	325	ARG	NE-CZ-NH1	15.23	127.91	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	PCA	Mainchain
1	B	1	PCA	Mainchain
1	C	1	PCA	Mainchain
1	D	1	PCA	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	3022	0	2905	36	1
1	B	3022	0	2905	42	1
1	C	3022	0	2905	49	0
1	D	3022	0	2905	47	0
2	E	23	0	21	1	0
2	F	23	0	21	1	0
2	G	23	0	21	1	0
2	H	23	0	21	2	0
3	A	28	0	26	0	0
3	B	28	0	26	0	0
3	C	28	0	26	0	0
3	D	28	0	26	0	0
4	A	277	0	0	2	0
4	B	268	0	0	4	0
4	C	268	0	0	4	0
4	D	288	0	0	5	0
All	All	13393	0	11808	173	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 7.

The worst 5 of 173 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:203:ALA:HB2	1:B:208:THR:HG23	1.42	1.00
1:C:203:ALA:HB2	1:C:208:THR:HG23	1.47	0.95
1:D:203:ALA:HB2	1:D:208:THR:HG23	1.48	0.94
1:C:5:LYS:H	1:C:5:LYS:HE2	1.34	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:LYS:H	1:B:5:LYS:HE2	1.35	0.91

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:A:290:ASP:O	1:B:321:ASN:ND2[2_656]	2.09	0.11

## 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	396/411 (96%)	381 (96%)	15 (4%)	0	100	100
1	B	396/411 (96%)	381 (96%)	15 (4%)	0	100	100
1	C	396/411 (96%)	382 (96%)	14 (4%)	0	100	100
1	D	396/411 (96%)	382 (96%)	14 (4%)	0	100	100
All	All	1584/1644 (96%)	1526 (96%)	58 (4%)	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	323/333 (97%)	303 (94%)	20 (6%)	18	25
1	B	323/333 (97%)	302 (94%)	21 (6%)	17	23
1	C	323/333 (97%)	300 (93%)	23 (7%)	14	19
1	D	323/333 (97%)	303 (94%)	20 (6%)	18	25
All	All	1292/1332 (97%)	1208 (94%)	84 (6%)	17	23

5 of 84 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	276	LYS
1	D	102	LEU
1	C	302	SER
1	D	20	LYS
1	D	162	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	96	GLN
1	C	277	GLN
1	D	96	GLN
1	B	96	GLN
1	A	96	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	PCA	A	1	1	7,8,9	1.25	1 (14%)	9,10,12	1.84	4 (44%)
1	PCA	C	1	1	7,8,9	1.52	1 (14%)	9,10,12	2.64	5 (55%)
1	PCA	D	1	1	7,8,9	1.61	1 (14%)	9,10,12	2.24	4 (44%)
1	PCA	B	1	1	7,8,9	1.21	0	9,10,12	1.70	3 (33%)

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
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1
1	PCA	C	1	1	-	0/0/11/13	0/1/1/1
1	PCA	D	1	1	-	0/0/11/13	0/1/1/1
1	PCA	B	1	1	-	0/0/11/13	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	1	PCA	CG-CD	2.56	1.57	1.50
1	C	1	PCA	O-C	2.10	1.28	1.19
1	A	1	PCA	O-C	2.10	1.28	1.19

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1	PCA	OE-CD-CG	-5.71	116.81	126.76
1	D	1	PCA	OE-CD-CG	-4.03	119.73	126.76
1	C	1	PCA	O-C-CA	-3.48	115.66	124.78
1	D	1	PCA	CB-CA-C	-3.34	108.10	112.70
1	A	1	PCA	OE-CD-CG	-3.20	121.18	126.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates i

8 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	BGC	E	1	2	12,12,12	1.01	0	17,17,17	1.80	4 (23%)
2	BGC	E	2	2	11,11,12	1.40	1 (9%)	15,15,17	1.18	2 (13%)
2	BGC	F	1	2	12,12,12	1.27	2 (16%)	17,17,17	1.71	4 (23%)
2	BGC	F	2	2	11,11,12	1.23	1 (9%)	15,15,17	1.39	2 (13%)
2	BGC	G	1	2	12,12,12	1.22	0	17,17,17	1.72	2 (11%)
2	BGC	G	2	2	11,11,12	1.35	1 (9%)	15,15,17	1.30	1 (6%)
2	BGC	H	1	2	12,12,12	1.17	0	17,17,17	1.50	3 (17%)
2	BGC	H	2	2	11,11,12	1.41	2 (18%)	15,15,17	1.63	3 (20%)

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	BGC	E	1	2	-	0/2/22/22	0/1/1/1
2	BGC	E	2	2	-	1/2/19/22	0/1/1/1
2	BGC	F	1	2	-	0/2/22/22	0/1/1/1
2	BGC	F	2	2	-	1/2/19/22	0/1/1/1
2	BGC	G	1	2	-	0/2/22/22	0/1/1/1
2	BGC	G	2	2	-	1/2/19/22	0/1/1/1
2	BGC	H	1	2	-	0/2/22/22	0/1/1/1
2	BGC	H	2	2	-	1/2/19/22	0/1/1/1

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	2	BGC	O5-C5	3.43	1.50	1.43
2	G	2	BGC	O5-C5	3.02	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	2	BGC	C2-C3	2.62	1.56	1.52
2	F	2	BGC	O5-C5	2.60	1.48	1.43
2	H	2	BGC	O5-C5	2.48	1.48	1.43

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	BGC	O3-C3-C2	-4.24	100.55	110.35
2	G	1	BGC	O3-C3-C2	-4.19	100.66	110.35
2	F	1	BGC	O3-C3-C2	-4.15	100.76	110.35
2	H	2	BGC	O2-C2-C1	-4.00	100.96	109.15
2	F	2	BGC	O2-C2-C1	-3.37	102.26	109.15

There are no chirality outliers.

All (4) torsion outliers are listed below:

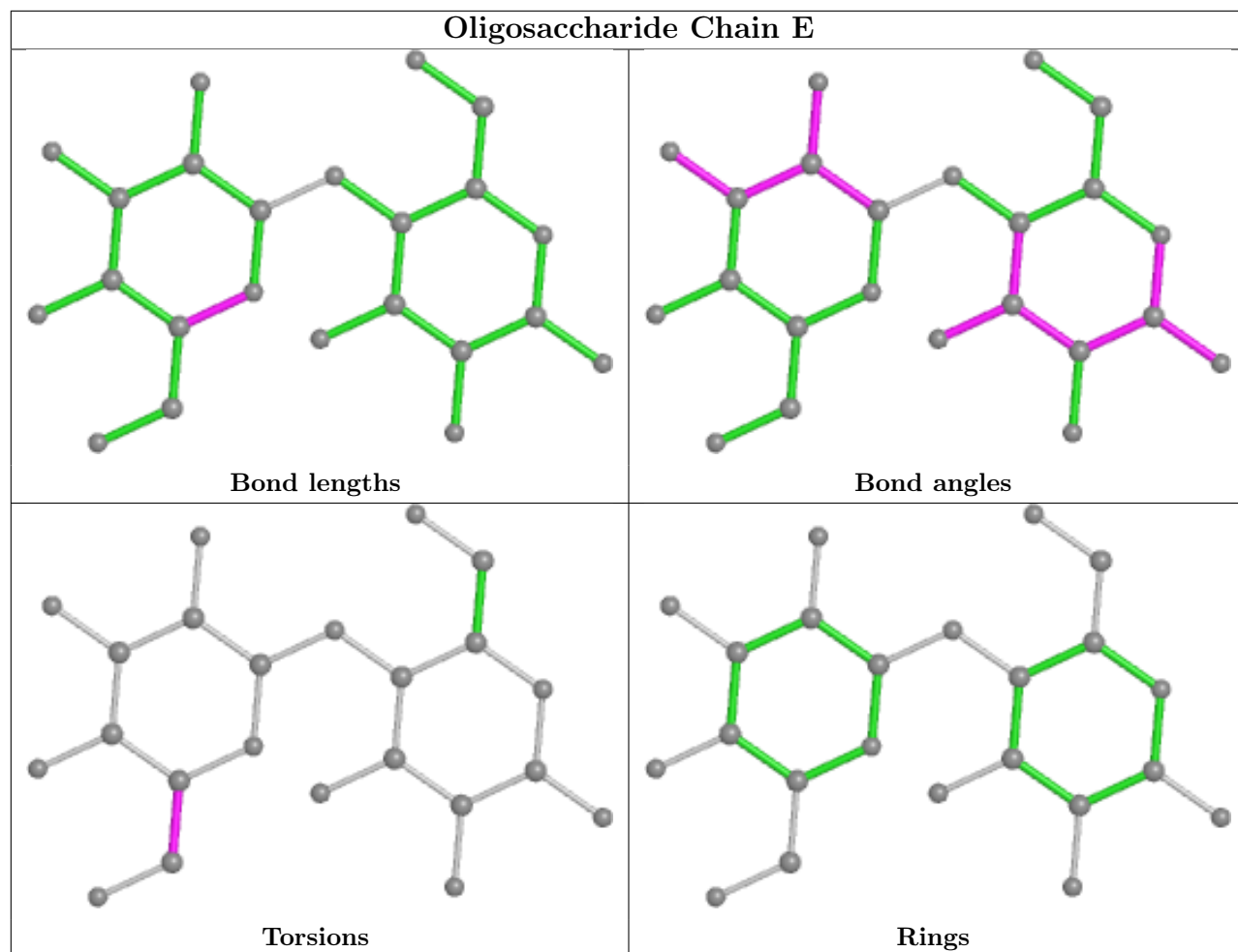
Mol	Chain	Res	Type	Atoms
2	H	2	BGC	C4-C5-C6-O6
2	F	2	BGC	C4-C5-C6-O6
2	E	2	BGC	C4-C5-C6-O6
2	G	2	BGC	C4-C5-C6-O6

There are no ring outliers.

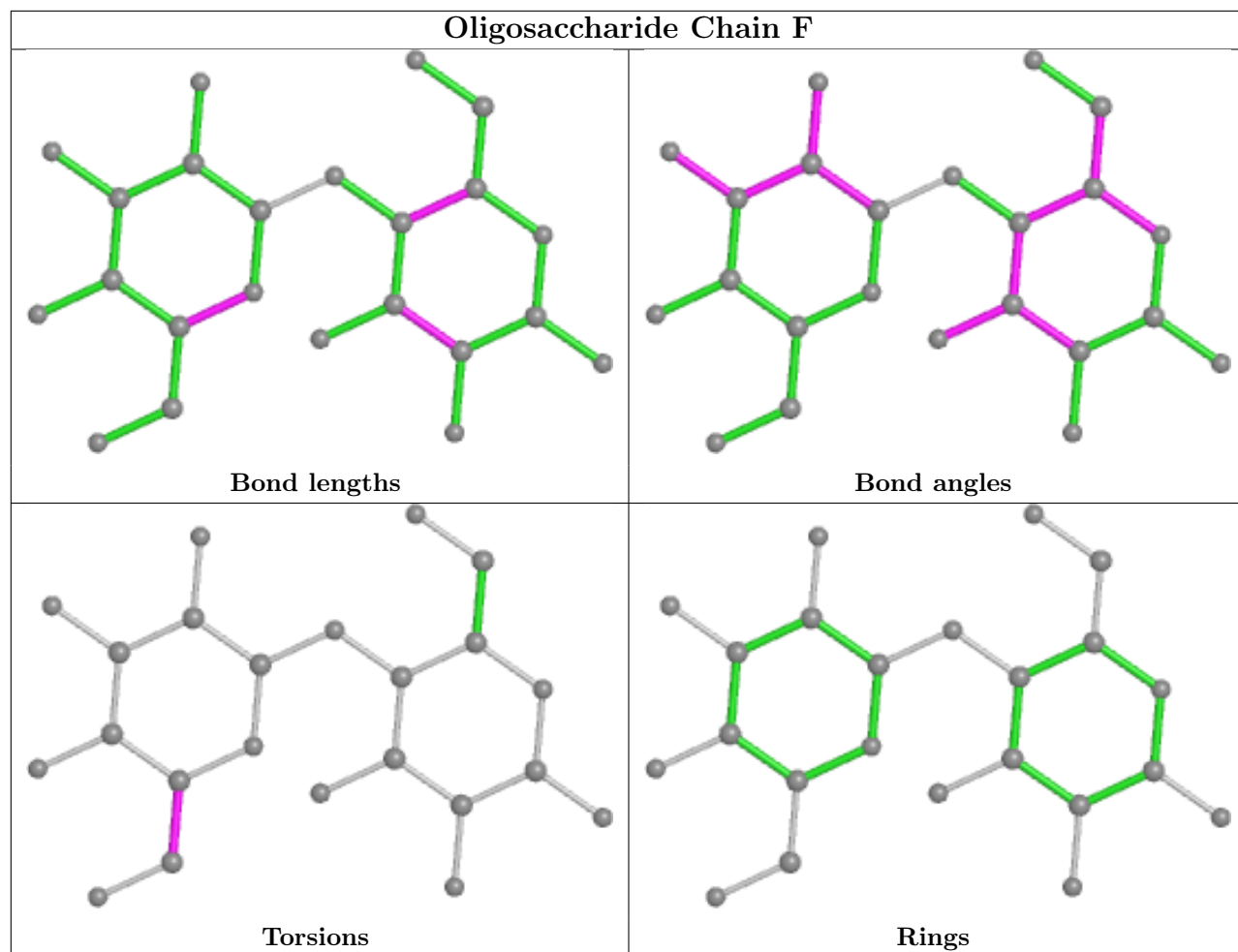
4 monomers are involved in 5 short contacts:

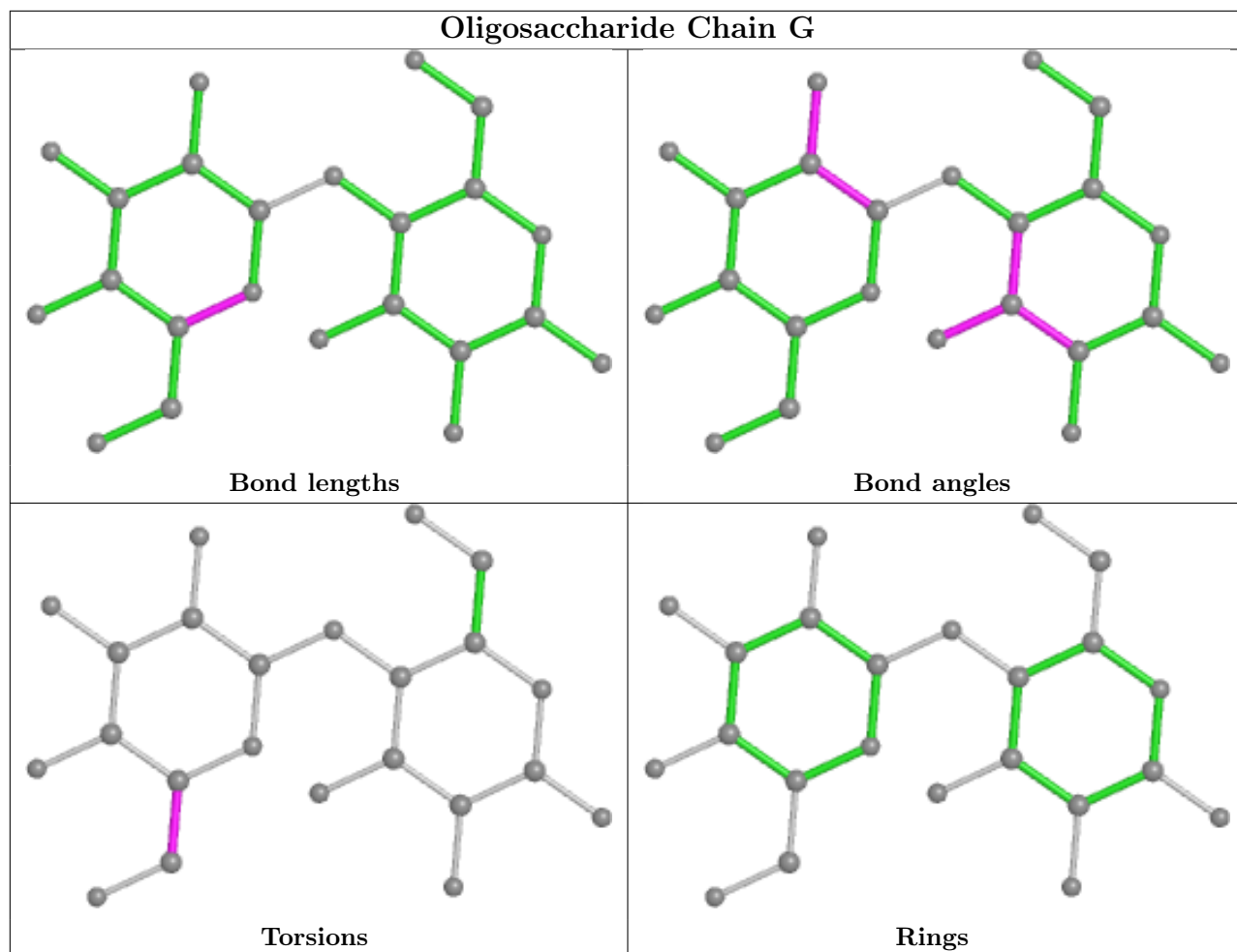
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	1	BGC	2	0
2	E	1	BGC	1	0
2	G	1	BGC	1	0
2	F	1	BGC	1	0

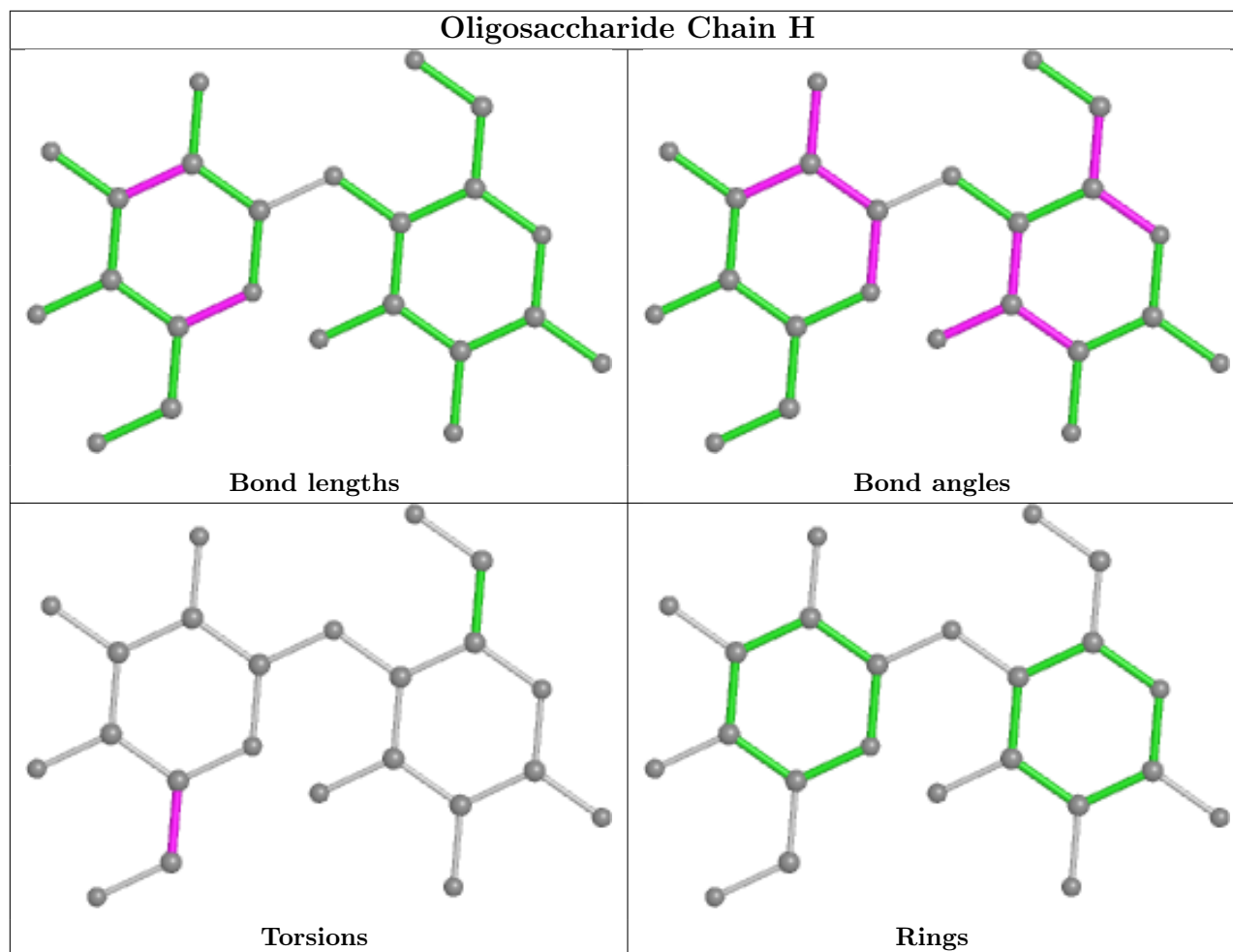
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











## 5.6 Ligand geometry [i](#)

8 ligands 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$
3	NAG	A	413	1	14,14,15	1.15	1 (7%)	17,19,21	1.53	3 (17%)
3	NAG	B	413	1	14,14,15	1.25	1 (7%)	17,19,21	1.98	7 (41%)
3	NAG	C	413	1	14,14,15	1.18	1 (7%)	17,19,21	1.57	2 (11%)
3	NAG	A	412	1	14,14,15	1.29	1 (7%)	17,19,21	1.55	4 (23%)
3	NAG	C	412	1	14,14,15	1.24	1 (7%)	17,19,21	2.18	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	B	412	1	14,14,15	1.27	1 (7%)	17,19,21	1.63	2 (11%)
3	NAG	D	413	1	14,14,15	1.22	1 (7%)	17,19,21	1.88	5 (29%)
3	NAG	D	412	1	14,14,15	1.36	1 (7%)	17,19,21	1.78	5 (29%)

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
3	NAG	A	413	1	-	2/6/23/26	0/1/1/1
3	NAG	B	413	1	-	0/6/23/26	0/1/1/1
3	NAG	C	413	1	-	2/6/23/26	0/1/1/1
3	NAG	A	412	1	-	0/6/23/26	0/1/1/1
3	NAG	C	412	1	-	2/6/23/26	0/1/1/1
3	NAG	B	412	1	-	0/6/23/26	0/1/1/1
3	NAG	D	413	1	-	2/6/23/26	0/1/1/1
3	NAG	D	412	1	-	2/6/23/26	0/1/1/1

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	413	NAG	O7-C7	-3.91	1.14	1.23
3	D	412	NAG	O7-C7	-3.73	1.14	1.23
3	A	412	NAG	O7-C7	-3.66	1.15	1.23
3	B	413	NAG	O7-C7	-3.52	1.15	1.23
3	B	412	NAG	O7-C7	-3.51	1.15	1.23

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	412	NAG	C1-O5-C5	-7.11	102.56	112.19
3	B	412	NAG	C1-O5-C5	-4.75	105.76	112.19
3	C	413	NAG	C1-O5-C5	4.46	118.24	112.19
3	D	413	NAG	C2-N2-C7	-4.24	116.86	122.90
3	D	412	NAG	O5-C1-C2	-4.19	104.67	111.29

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	413	NAG	O5-C5-C6-O6
3	A	413	NAG	C8-C7-N2-C2
3	A	413	NAG	O7-C7-N2-C2
3	C	412	NAG	C4-C5-C6-O6
3	D	413	NAG	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, 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	A	397/411 (96%)	-0.18	9 (2%) 60 67	13, 26, 43, 52	0
1	B	397/411 (96%)	0.05	18 (4%) 33 40	14, 27, 44, 51	0
1	C	397/411 (96%)	-0.07	8 (2%) 65 71	12, 27, 44, 53	0
1	D	397/411 (96%)	-0.21	6 (1%) 73 79	14, 25, 42, 51	0
All	All	1588/1644 (96%)	-0.10	41 (2%) 56 63	12, 26, 43, 53	0

The worst 5 of 41 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	159	ARG	3.8
1	C	201	TRP	3.0
1	B	368	THR	2.9
1	C	276	LYS	2.9
1	B	203	ALA	2.9

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 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.

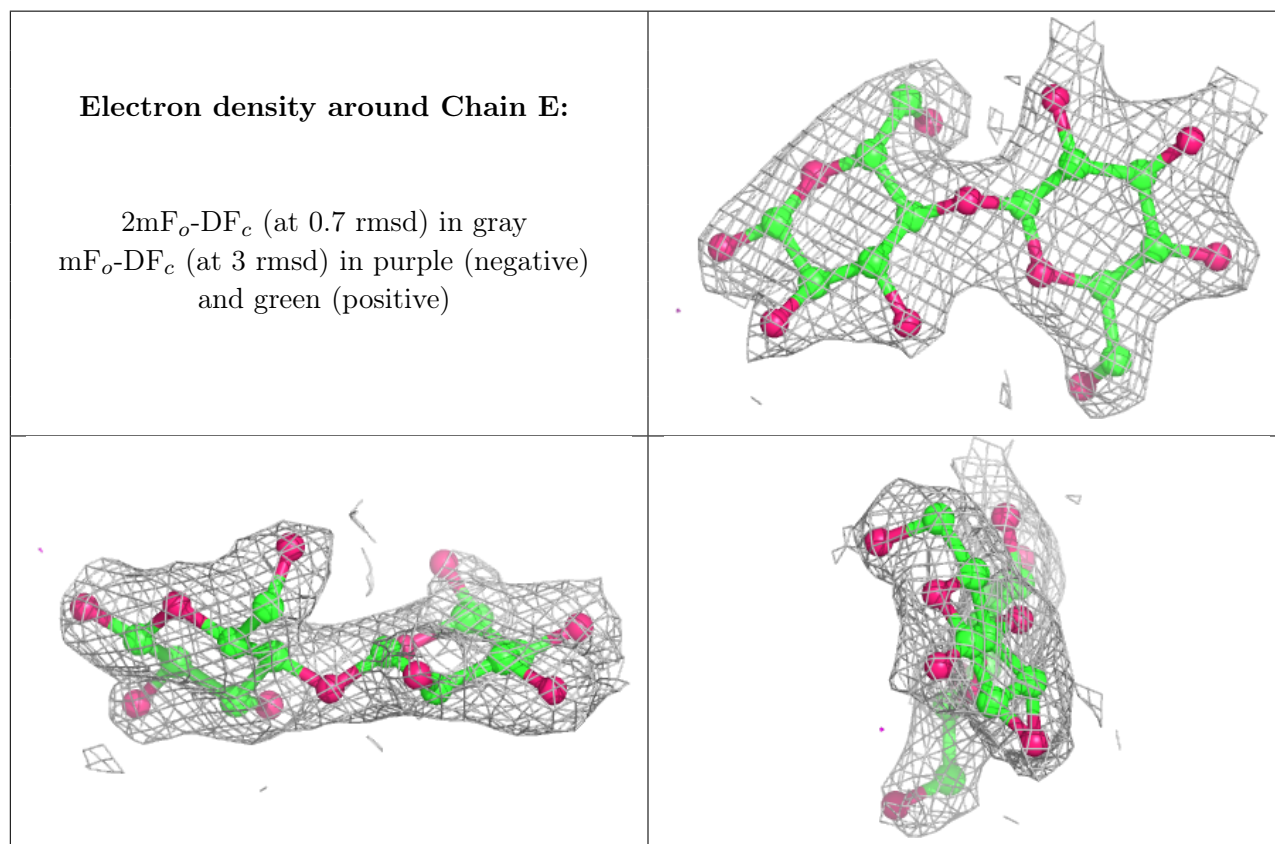
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	PCA	C	1	8/9	0.94	0.12	27,29,31,33	0
1	PCA	B	1	8/9	0.96	0.09	24,27,29,31	0
1	PCA	A	1	8/9	0.96	0.09	23,26,29,31	0
1	PCA	D	1	8/9	0.96	0.09	23,26,27,28	0

### 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, 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.

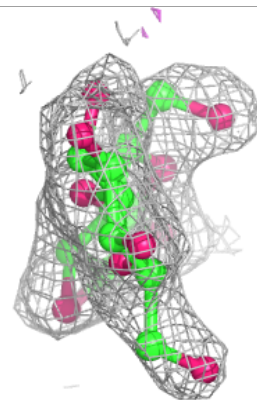
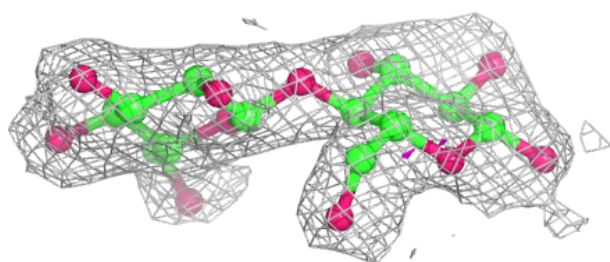
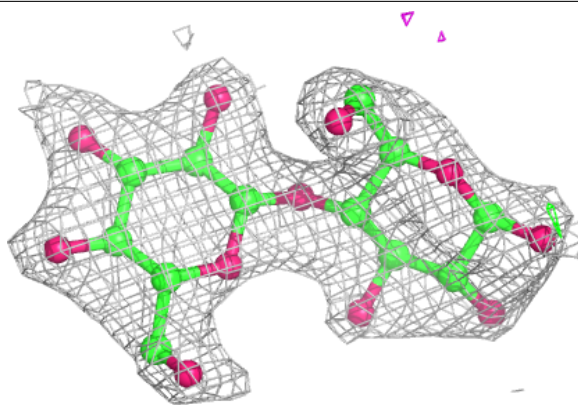
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	BGC	G	1	12/12	0.88	0.20	32,35,38,39	0
2	BGC	H	2	11/12	0.90	0.16	26,33,39,43	0
2	BGC	G	2	11/12	0.91	0.16	28,33,40,43	0
2	BGC	F	2	11/12	0.91	0.17	28,32,39,43	0
2	BGC	E	2	11/12	0.92	0.14	28,33,39,43	0
2	BGC	F	1	12/12	0.92	0.15	34,35,38,38	0
2	BGC	H	1	12/12	0.93	0.18	33,35,38,39	0
2	BGC	E	1	12/12	0.93	0.19	33,35,37,39	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.

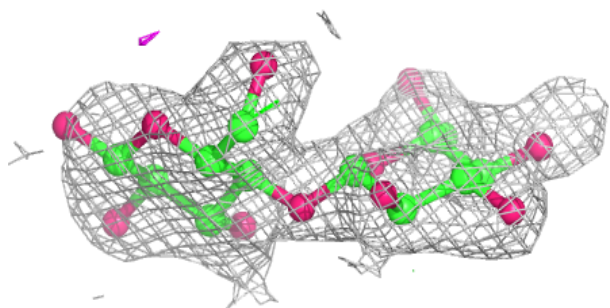
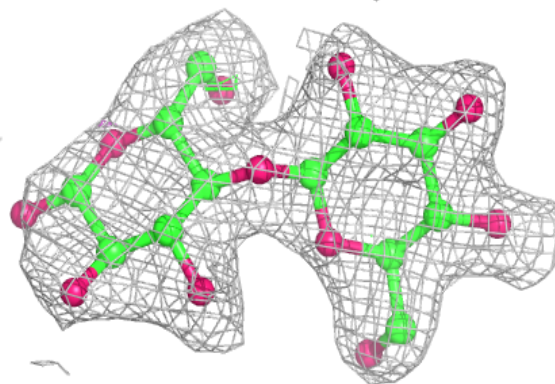


**Electron density around Chain F:**

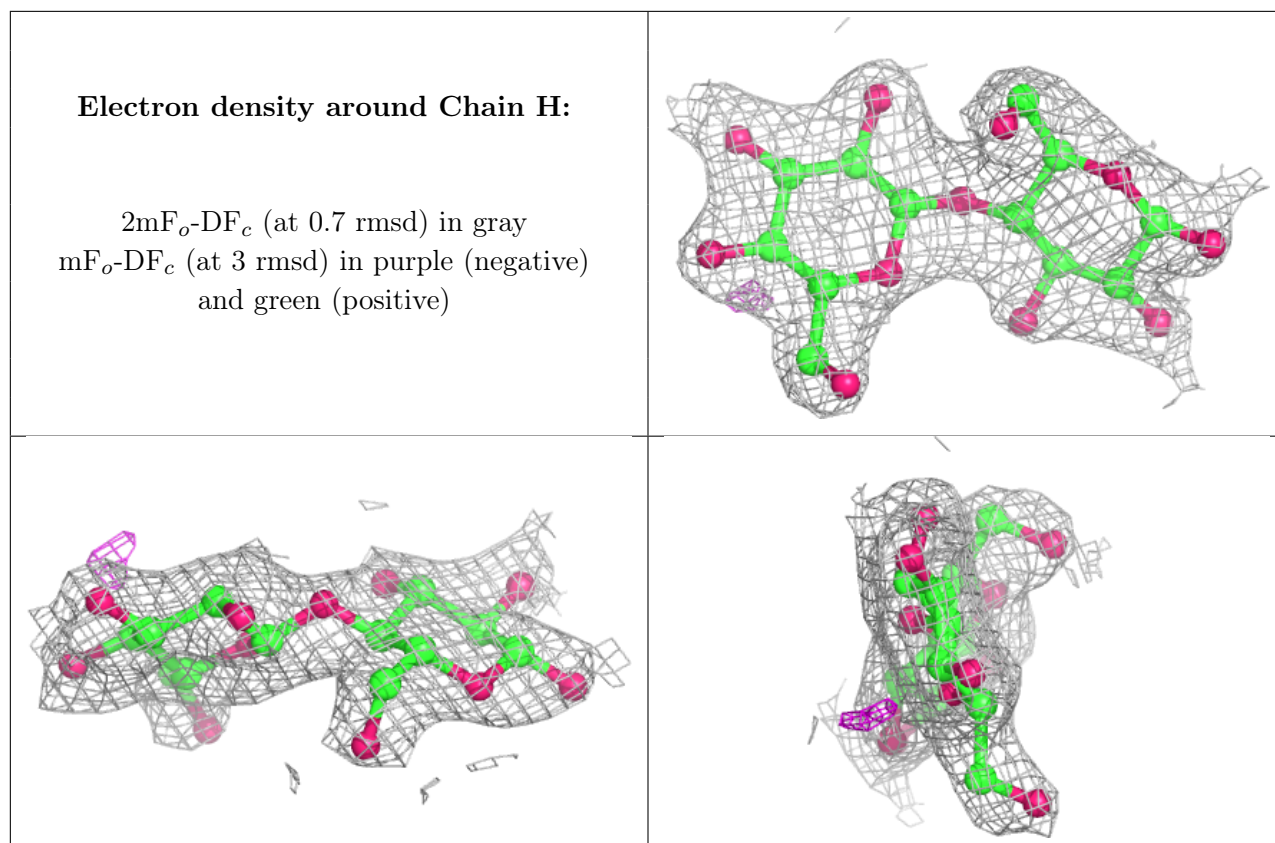
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around Chain 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, 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	NAG	D	413	14/15	0.77	0.27	57,59,62,62	0
3	NAG	B	413	14/15	0.81	0.25	59,61,63,63	0
3	NAG	A	413	14/15	0.83	0.23	56,59,62,62	0
3	NAG	A	412	14/15	0.87	0.25	53,55,56,56	0
3	NAG	B	412	14/15	0.87	0.23	52,54,55,56	0
3	NAG	C	413	14/15	0.88	0.18	56,58,59,61	0
3	NAG	D	412	14/15	0.88	0.19	52,54,57,57	0
3	NAG	C	412	14/15	0.88	0.16	52,54,55,55	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.