



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 3, 2023 – 04:57 PM EDT

PDB ID : 6OZ4  
Title : Crystal structure of broadly neutralizing antibody N49P6 Fab in complex with HIV-1 BG505 SOSIP.664 Env trimer ectodomain.  
Authors : Tolbert, W.D.; Pazgier, M.  
Deposited on : 2019-05-15  
Resolution : 4.05 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 : 2.35.1  
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) : 2.35.1

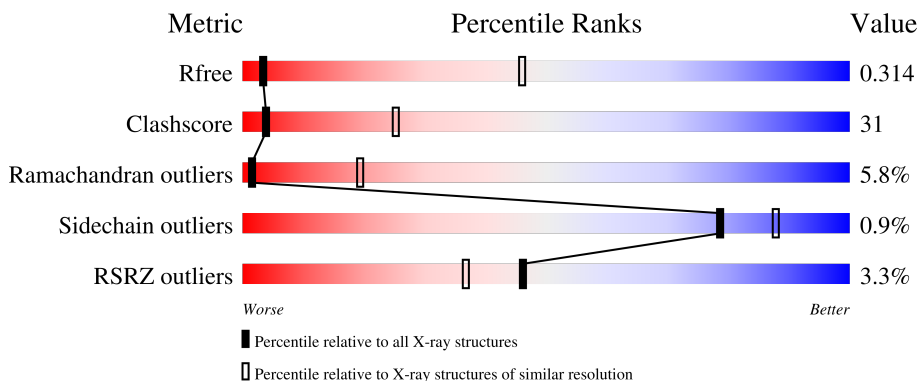
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 4.05 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.



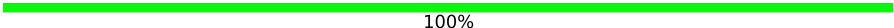


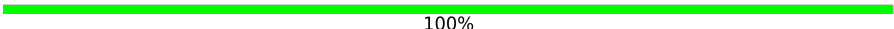

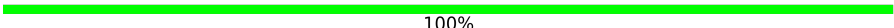
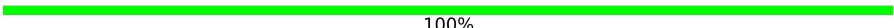
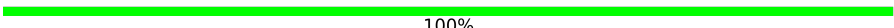
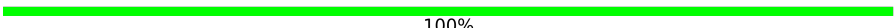
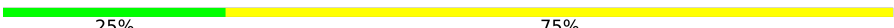
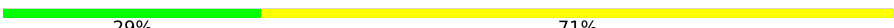
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1127 (4.42-3.70)
Clashscore	141614	1033 (4.40-3.72)
Ramachandran outliers	138981	1145 (4.42-3.70)
Sidechain outliers	138945	1133 (4.42-3.70)
RSRZ outliers	127900	1005 (4.44-3.68)

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	G	481	
2	B	153	
3	H	229	
4	L	205	
5	A	3	

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Mol	Chain	Length	Quality of chain
5	F	3	 100%
5	M	3	 33% 67%
6	C	2	 50% 50%
6	E	2	 100%
6	I	2	 50% 50%
6	J	2	 100%
6	N	2	 100%
6	O	2	 100%
6	P	2	 100%
7	D	4	 25% 75%
8	K	7	 29% 71%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	BMA	A	3	-	-	-	X
6	NAG	C	2	-	-	-	X
6	NAG	E	2	-	-	-	X
6	NAG	I	1	-	-	-	X
7	BMA	D	3	-	-	-	X
7	MAN	D	4	-	-	-	X
9	NAG	G	603	-	-	-	X

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 8269 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 Envelope glycoprotein gp160.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	G	449	3536	2219	625	664	28	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	332	ASN	THR	engineered mutation	UNP Q2N0S6
G	501	CYS	ALA	engineered mutation	UNP Q2N0S6
G	509	ARG	-	expression tag	UNP Q2N0S6
G	510	ARG	-	expression tag	UNP Q2N0S6
G	511	ARG	-	expression tag	UNP Q2N0S6
G	512	ARG	-	expression tag	UNP Q2N0S6
G	513	ARG	-	expression tag	UNP Q2N0S6

- Molecule 2 is a protein called Envelope glycoprotein gp41.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	122	971	611	168	186	6	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	559	PRO	ILE	engineered mutation	UNP Q2N0S6
B	605	CYS	THR	engineered mutation	UNP Q2N0S6

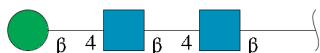
- Molecule 3 is a protein called N49P6 antibody Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	H	225	1722	1082	303	329	8	0	0	0

- Molecule 4 is a protein called N49P6 antibody light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	L	200	1496	936	254	301	5	0	0	0

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	A	3	39	22	2	15	0	0	0
5	F	3	39	22	2	15	0	0	0
5	M	3	39	22	2	15	0	0	0

- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



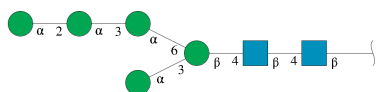
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
6	C	2	28	16	2	10	0	0	0
6	E	2	28	16	2	10	0	0	0
6	I	2	28	16	2	10	0	0	0
6	J	2	28	16	2	10	0	0	0
6	N	2	28	16	2	10	0	0	0
6	O	2	28	16	2	10	0	0	0
6	P	2	28	16	2	10	0	0	0

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



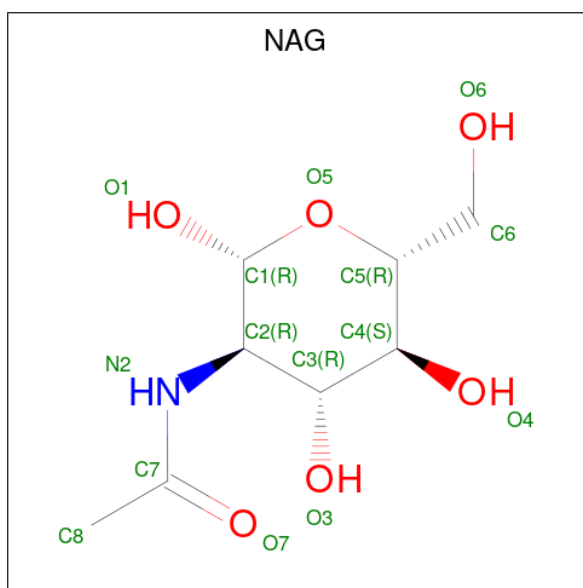
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
7	D	4	50	28	2	20	0	0	0

- Molecule 8 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
8	K	7	83	46	2	35	0	0	0

- Molecule 9 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>).

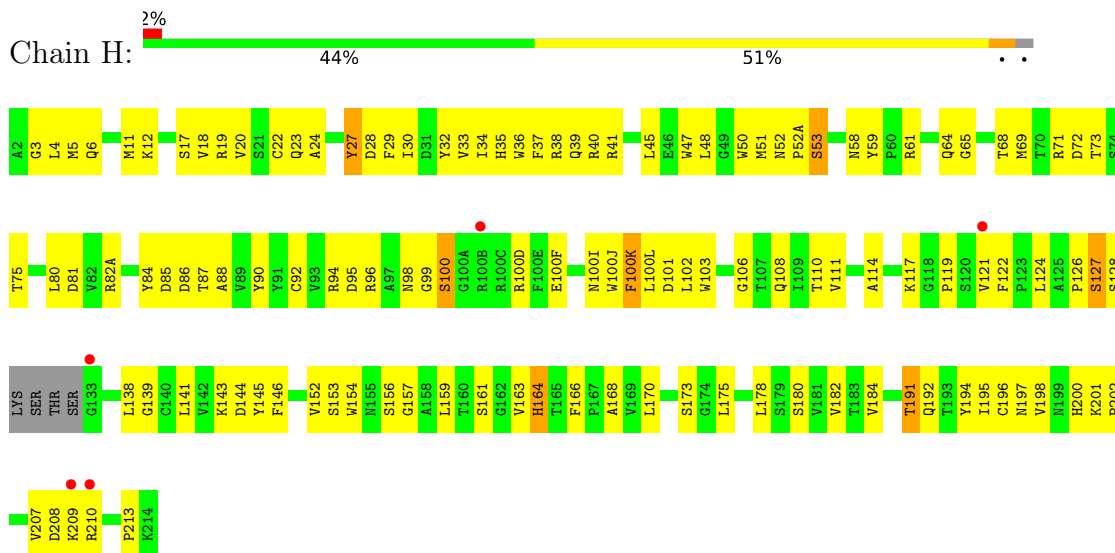


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	G	1	Total 14	C 8	N 1	O 5	0	0
9	G	1	Total 14	C 8	N 1	O 5	0	0
9	G	1	Total 14	C 8	N 1	O 5	0	0
9	G	1	Total 14	C 8	N 1	O 5	0	0
9	B	1	Total 14	C 8	N 1	O 5	0	0
9	B	1	Total 14	C 8	N 1	O 5	0	0
9	B	1	Total 14	C 8	N 1	O 5	0	0

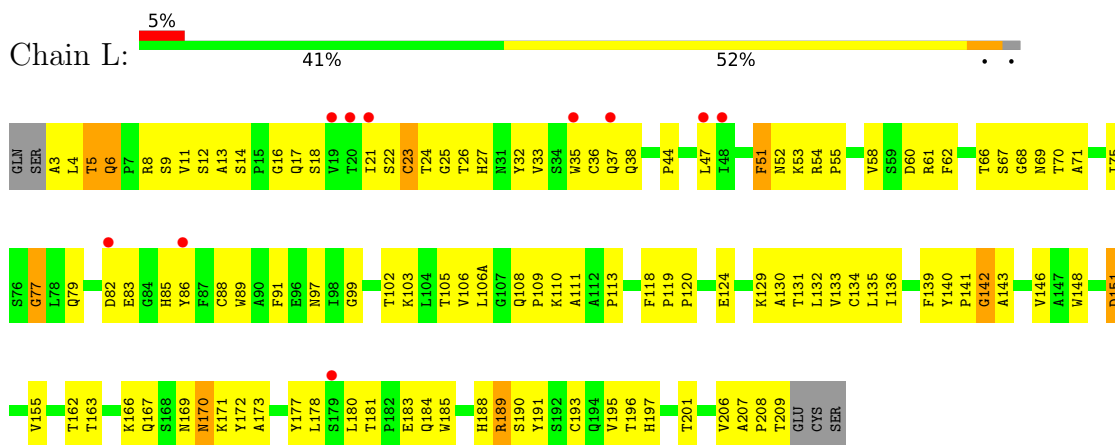




- Molecule 3: N49P6 antibody Fab heavy chain



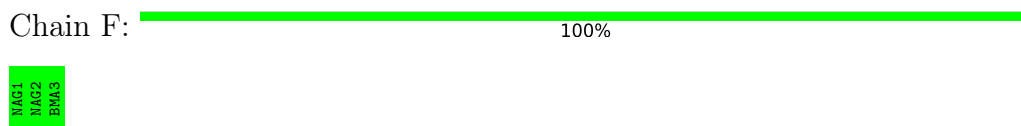
- Molecule 4: N49P6 antibody light chain



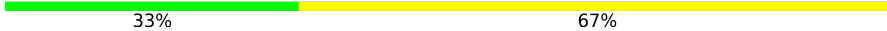
- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose




- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  33% 67%

  
MAG1  
MAG2  
MAG3

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  50% 50%

  
MAG1  
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  100%

  
MAG1  
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  50% 50%

  
MAG1  
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  100%

  
MAG1  
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  100%

  
MAG1  
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  100%


  
MAG1  
MAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:  100%

MAG1  
MAG2

- Molecule 7: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  25% 75%

MAG1  
MAG2  
BMA3  
MAN4

- Molecule 8: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  29% 71%

MAG1  
MAG2  
BMA3  
MAN4  
MAN5  
MAN6  
MAN7

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	164.70Å 164.70Å 164.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.95 – 4.05 49.66 – 4.03	Depositor EDS
% Data completeness (in resolution range)	99.4 (39.95-4.05) 99.7 (49.66-4.03)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.15 (at 4.00Å)	Xtrriage
Refinement program	PHENIX (1.15.2_3472: ???), REFMAC	Depositor
R, $R_{free}$	0.255 , 0.315 0.254 , 0.314	Depositor DCC
$R_{free}$ test set	580 reflections (4.62%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	208.8	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 245.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.072 for l,-k,h	Xtrriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	8269	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	225.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.99% 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: BMA, NAG, MAN

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	G	0.42	0/3610	0.68	0/4901
2	B	0.38	0/988	0.63	0/1340
3	H	0.44	0/1765	0.70	0/2401
4	L	0.40	0/1534	0.67	0/2094
All	All	0.42	0/7897	0.68	0/10736

There are no bond length outliers.

There are no bond angle outliers.

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	G	3536	0	3471	251	0
2	B	971	0	945	65	0
3	H	1722	0	1664	105	0
4	L	1496	0	1447	114	0
5	A	39	0	34	0	0
5	F	39	0	34	0	0
5	M	39	0	34	2	0
6	C	28	0	25	2	0
6	E	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	I	28	0	25	0	0
6	J	28	0	25	0	0
6	N	28	0	25	0	0
6	O	28	0	25	0	0
6	P	28	0	25	0	0
7	D	50	0	43	1	0
8	K	83	0	70	2	0
9	B	42	0	39	2	0
9	G	56	0	52	4	0
All	All	8269	0	8008	499	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 31.

The worst 5 of 499 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:G:43:PRO:CG	2:B:628:TRP:CD1	2.09	1.33
1:G:43:PRO:HG2	2:B:628:TRP:CD1	1.72	1.19
1:G:101:VAL:HB	1:G:483:LEU:HD12	1.21	1.13
1:G:43:PRO:HG2	2:B:628:TRP:CG	1.84	1.10
1:G:184:ILE:HD11	1:G:190:GLU:HG2	1.27	1.09

There are no symmetry-related clashes.

## 5.3 Torsion angles [\(i\)](#)

### 5.3.1 Protein backbone [\(i\)](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	G	443/481 (92%)	351 (79%)	61 (14%)	31 (7%)	<b>1</b> <b>16</b>
2	B	118/153 (77%)	92 (78%)	21 (18%)	5 (4%)	<b>3</b> <b>25</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	H	221/229 (96%)	190 (86%)	22 (10%)	9 (4%)	3	25
4	L	198/205 (97%)	163 (82%)	23 (12%)	12 (6%)	1	18
All	All	980/1068 (92%)	796 (81%)	127 (13%)	57 (6%)	1	19

5 of 57 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	65	LYS
1	G	66	HIS
1	G	268	GLU
1	G	429	ARG
2	B	598	CYS

### 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	G	401/428 (94%)	399 (100%)	2 (0%)	88	93
2	B	105/129 (81%)	105 (100%)	0	100	100
3	H	188/192 (98%)	184 (98%)	4 (2%)	53	72
4	L	168/173 (97%)	166 (99%)	2 (1%)	71	83
All	All	862/922 (94%)	854 (99%)	8 (1%)	78	88

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	L	23	CYS
4	L	8	ARG
3	H	100(K)	PHE
3	H	100	SER
3	H	164	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
3	H	35	HIS
3	H	58	ASN
4	L	85	HIS
3	H	197	ASN
2	B	650	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](#)

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

### 5.5 Carbohydrates [i](#)

34 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	$\# Z  > 2$
5	NAG	A	1	5,1	14,14,15	0.40	0	17,19,21	0.37	0
5	NAG	A	2	5	14,14,15	0.49	0	17,19,21	0.68	1 (5%)
5	BMA	A	3	5	11,11,12	0.67	0	15,15,17	0.94	1 (6%)
6	NAG	C	1	1,6	14,14,15	0.25	0	17,19,21	0.45	0
6	NAG	C	2	6	14,14,15	0.35	0	17,19,21	1.39	2 (11%)
7	NAG	D	1	1,7	14,14,15	0.44	0	17,19,21	0.69	0
7	NAG	D	2	7	14,14,15	0.19	0	17,19,21	0.70	0
7	BMA	D	3	7	11,11,12	1.10	1 (9%)	15,15,17	1.11	1 (6%)
7	MAN	D	4	7	11,11,12	1.06	1 (9%)	15,15,17	1.46	3 (20%)
6	NAG	E	1	1,6	14,14,15	0.39	0	17,19,21	0.47	0
6	NAG	E	2	6	14,14,15	0.29	0	17,19,21	0.44	0
5	NAG	F	1	5,1	14,14,15	0.31	0	17,19,21	0.36	0
5	NAG	F	2	5	14,14,15	0.18	0	17,19,21	0.40	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	BMA	F	3	5	11,11,12	0.80	0	15,15,17	0.88	0
6	NAG	I	1	1,6	14,14,15	0.20	0	17,19,21	0.72	1 (5%)
6	NAG	I	2	6	14,14,15	0.30	0	17,19,21	0.41	0
6	NAG	J	1	1,6	14,14,15	0.22	0	17,19,21	0.45	0
6	NAG	J	2	6	14,14,15	0.32	0	17,19,21	0.51	0
8	NAG	K	1	8,1	14,14,15	0.46	0	17,19,21	0.50	0
8	NAG	K	2	8	14,14,15	0.26	0	17,19,21	0.59	0
8	BMA	K	3	8	11,11,12	0.70	0	15,15,17	0.94	0
8	MAN	K	4	8	11,11,12	0.72	0	15,15,17	1.31	2 (13%)
8	MAN	K	5	8	11,11,12	0.89	0	15,15,17	1.15	3 (20%)
8	MAN	K	6	8	11,11,12	0.82	0	15,15,17	1.29	2 (13%)
8	MAN	K	7	8	11,11,12	0.82	0	15,15,17	1.12	2 (13%)
5	NAG	M	1	5,1	14,14,15	0.31	0	17,19,21	0.54	0
5	NAG	M	2	5	14,14,15	0.42	0	17,19,21	0.65	0
5	BMA	M	3	5	11,11,12	0.92	1 (9%)	15,15,17	1.25	1 (6%)
6	NAG	N	1	1,6	14,14,15	0.43	0	17,19,21	0.75	0
6	NAG	N	2	6	14,14,15	0.29	0	17,19,21	0.41	0
6	NAG	O	1	1,6	14,14,15	0.36	0	17,19,21	0.68	0
6	NAG	O	2	6	14,14,15	0.27	0	17,19,21	0.49	0
6	NAG	P	1	1,6	14,14,15	0.45	0	17,19,21	0.56	0
6	NAG	P	2	6	14,14,15	0.24	0	17,19,21	0.51	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	A	2	5	-	1/6/23/26	0/1/1/1
5	BMA	A	3	5	-	1/2/19/22	0/1/1/1
6	NAG	C	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	C	2	6	-	5/6/23/26	0/1/1/1
7	NAG	D	1	1,7	-	3/6/23/26	0/1/1/1
7	NAG	D	2	7	-	3/6/23/26	0/1/1/1
7	BMA	D	3	7	-	2/2/19/22	0/1/1/1
7	MAN	D	4	7	-	0/2/19/22	0/1/1/1
6	NAG	E	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	E	2	6	-	3/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	F	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	F	2	5	-	1/6/23/26	0/1/1/1
5	BMA	F	3	5	-	2/2/19/22	0/1/1/1
6	NAG	I	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	I	2	6	-	1/6/23/26	0/1/1/1
6	NAG	J	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	J	2	6	-	2/6/23/26	0/1/1/1
8	NAG	K	1	8,1	-	0/6/23/26	0/1/1/1
8	NAG	K	2	8	-	3/6/23/26	0/1/1/1
8	BMA	K	3	8	-	0/2/19/22	0/1/1/1
8	MAN	K	4	8	-	1/2/19/22	0/1/1/1
8	MAN	K	5	8	-	0/2/19/22	0/1/1/1
8	MAN	K	6	8	-	2/2/19/22	1/1/1/1
8	MAN	K	7	8	-	0/2/19/22	0/1/1/1
5	NAG	M	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	M	2	5	-	3/6/23/26	0/1/1/1
5	BMA	M	3	5	-	0/2/19/22	0/1/1/1
6	NAG	N	1	1,6	-	3/6/23/26	0/1/1/1
6	NAG	N	2	6	-	2/6/23/26	0/1/1/1
6	NAG	O	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	O	2	6	-	2/6/23/26	0/1/1/1
6	NAG	P	1	1,6	-	4/6/23/26	0/1/1/1
6	NAG	P	2	6	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	3	BMA	O5-C1	-2.75	1.39	1.43
5	M	3	BMA	C1-C2	2.17	1.57	1.52
7	D	4	MAN	C2-C3	2.14	1.55	1.52

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	2	NAG	C2-N2-C7	4.56	129.39	122.90
8	K	6	MAN	C1-O5-C5	3.69	117.20	112.19
8	K	4	MAN	C1-O5-C5	3.04	116.32	112.19
8	K	7	MAN	C1-O5-C5	2.73	115.89	112.19
7	D	4	MAN	C1-O5-C5	2.64	115.76	112.19

There are no chirality outliers.

5 of 64 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	M	1	NAG	O5-C5-C6-O6
6	E	1	NAG	O5-C5-C6-O6
8	K	2	NAG	O5-C5-C6-O6
5	A	1	NAG	O5-C5-C6-O6
5	F	3	BMA	O5-C5-C6-O6

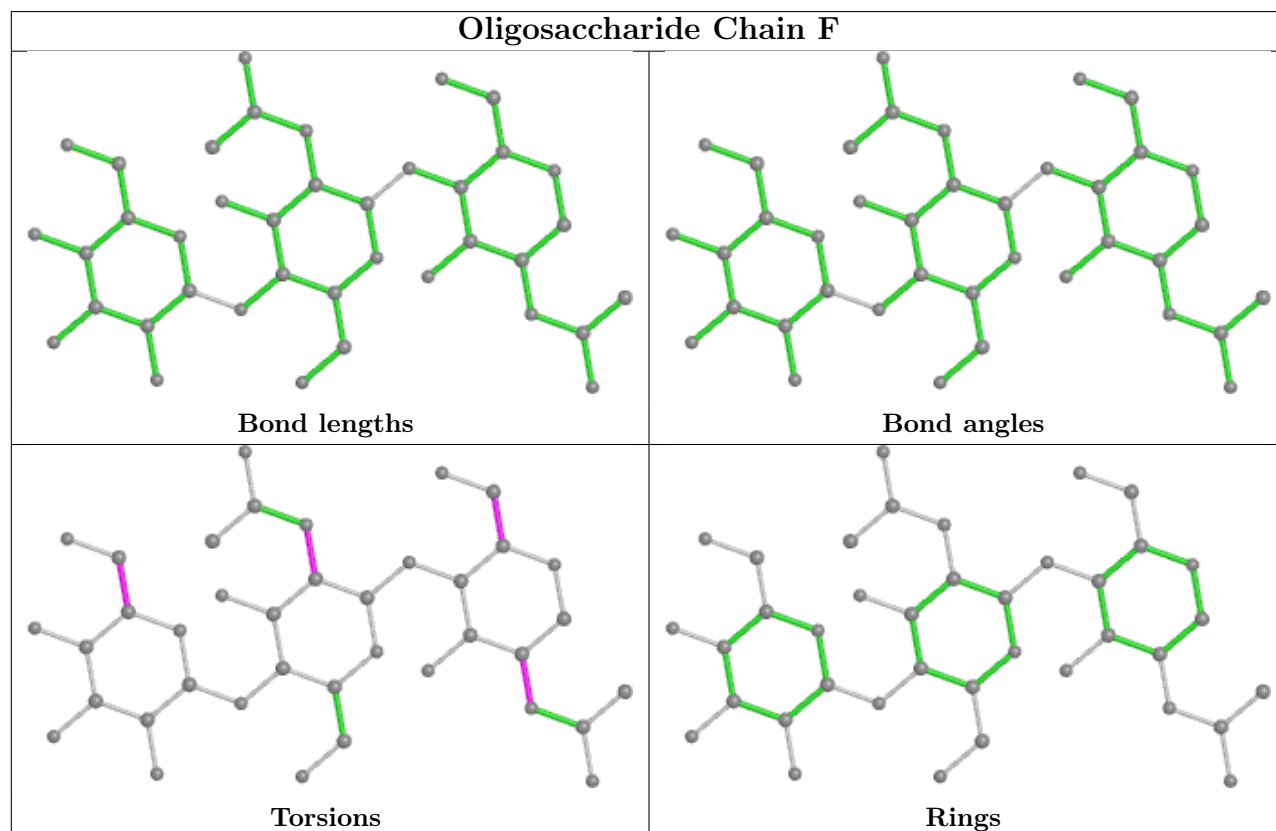
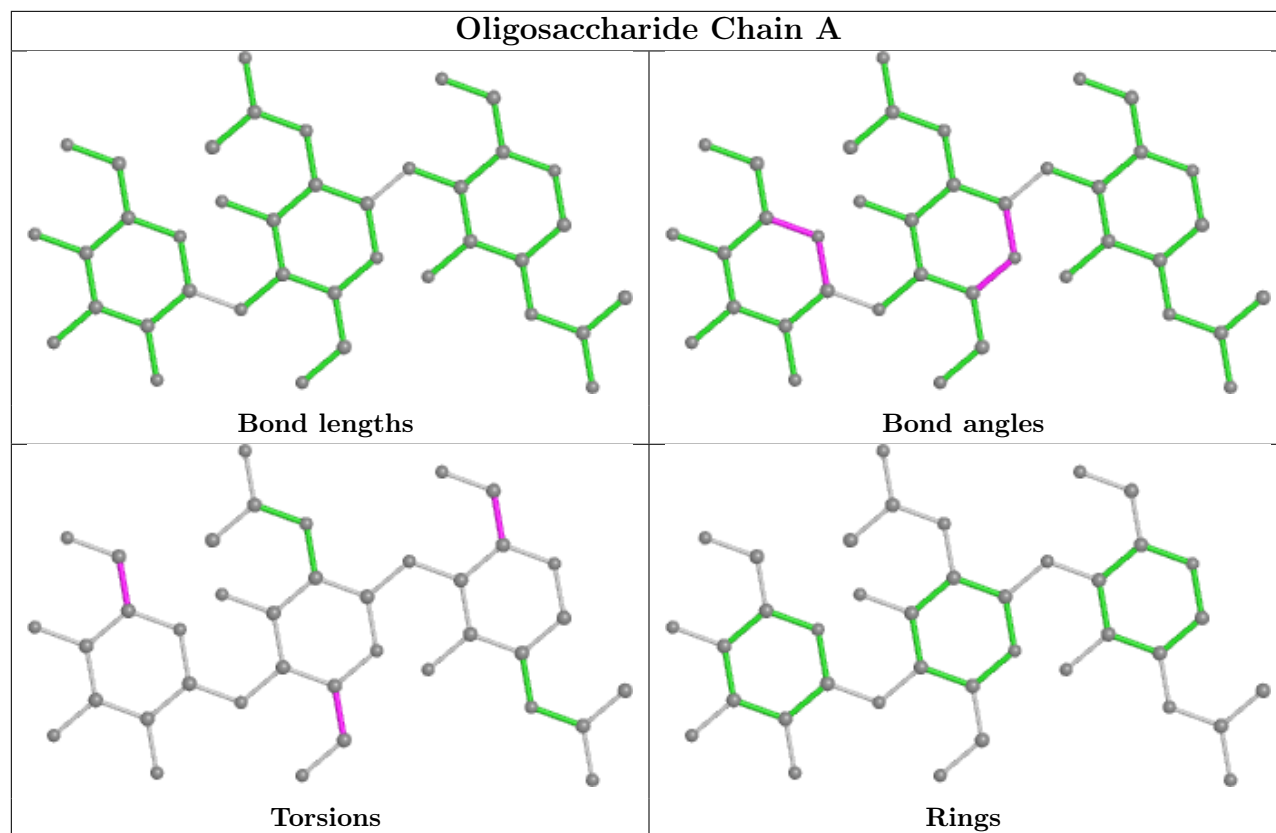
All (1) ring outliers are listed below:

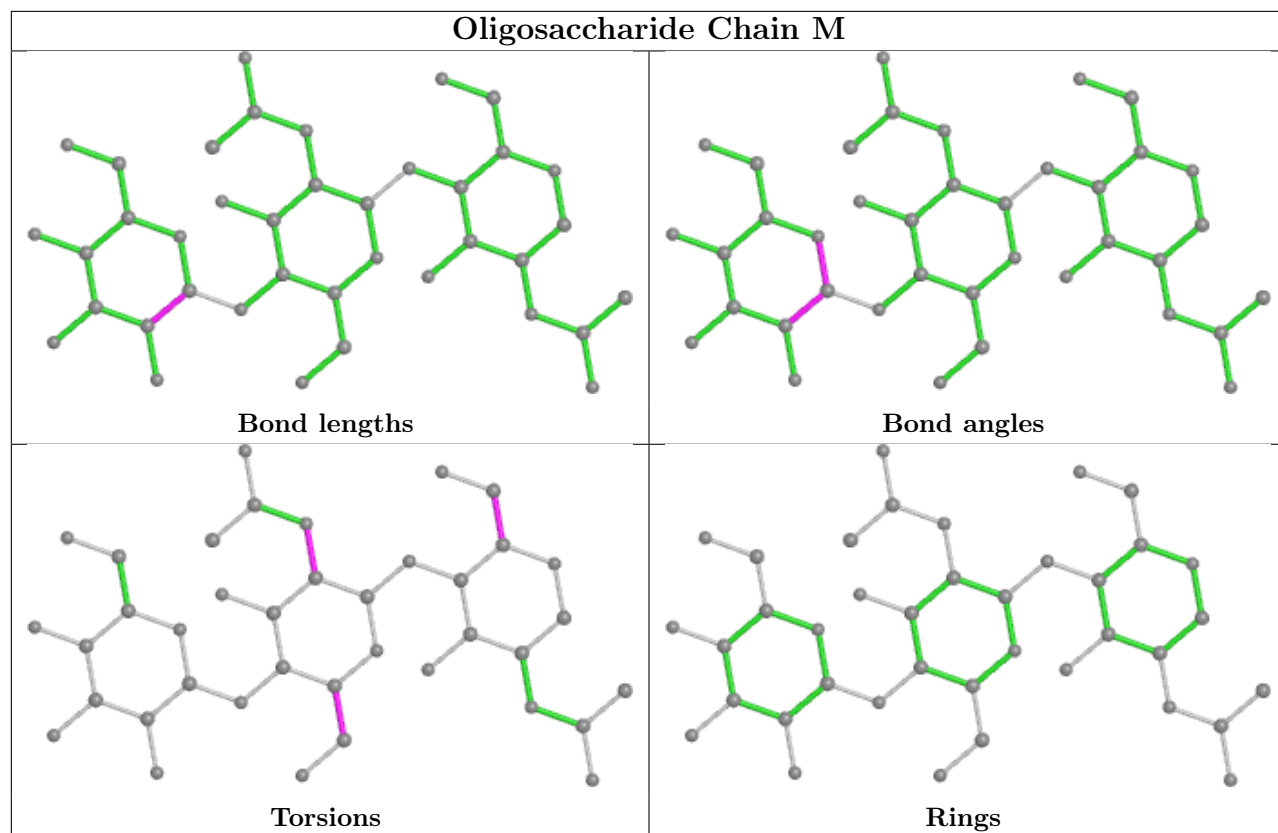
Mol	Chain	Res	Type	Atoms
8	K	6	MAN	C1-C2-C3-C4-C5-O5

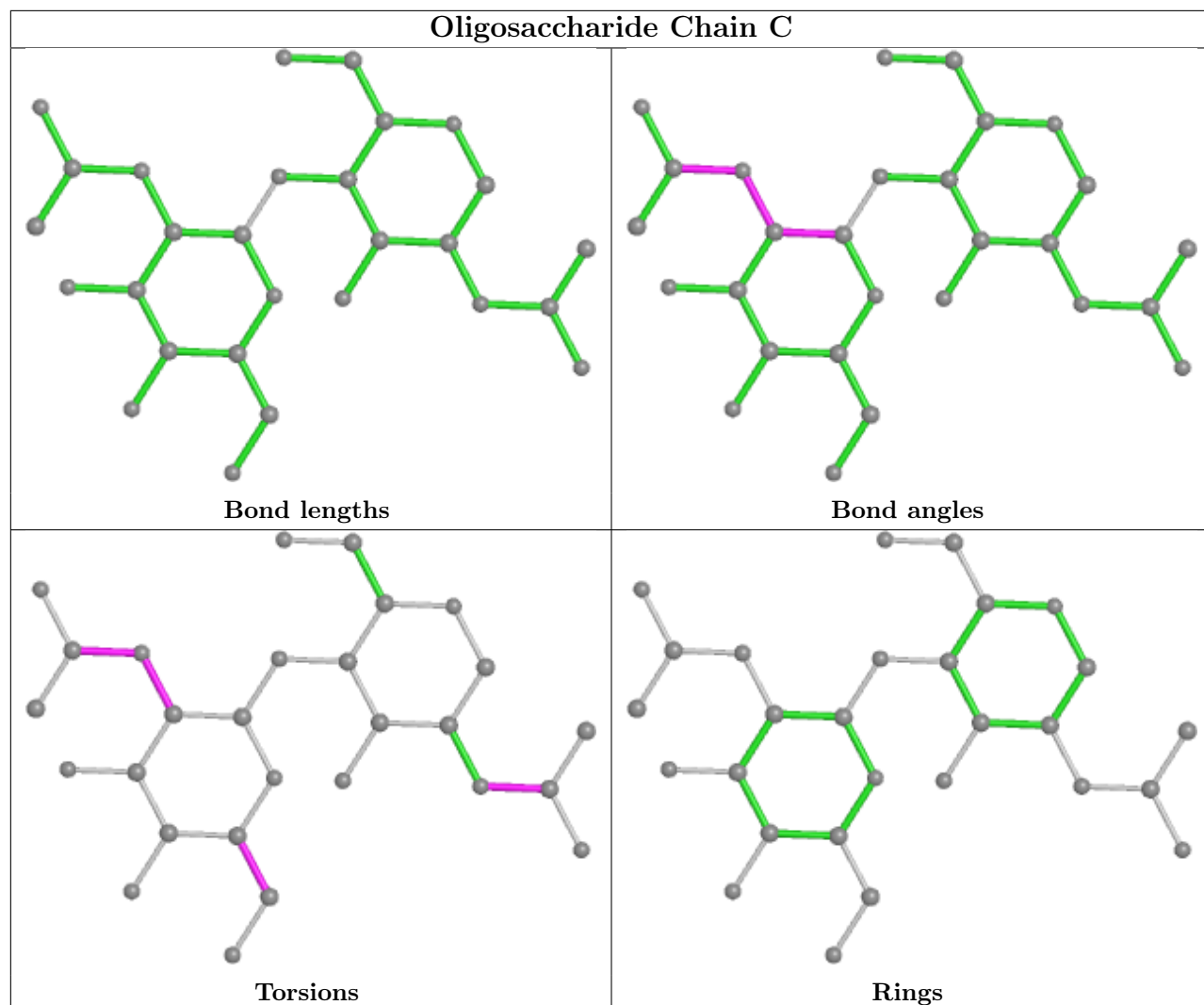
5 monomers are involved in 7 short contacts:

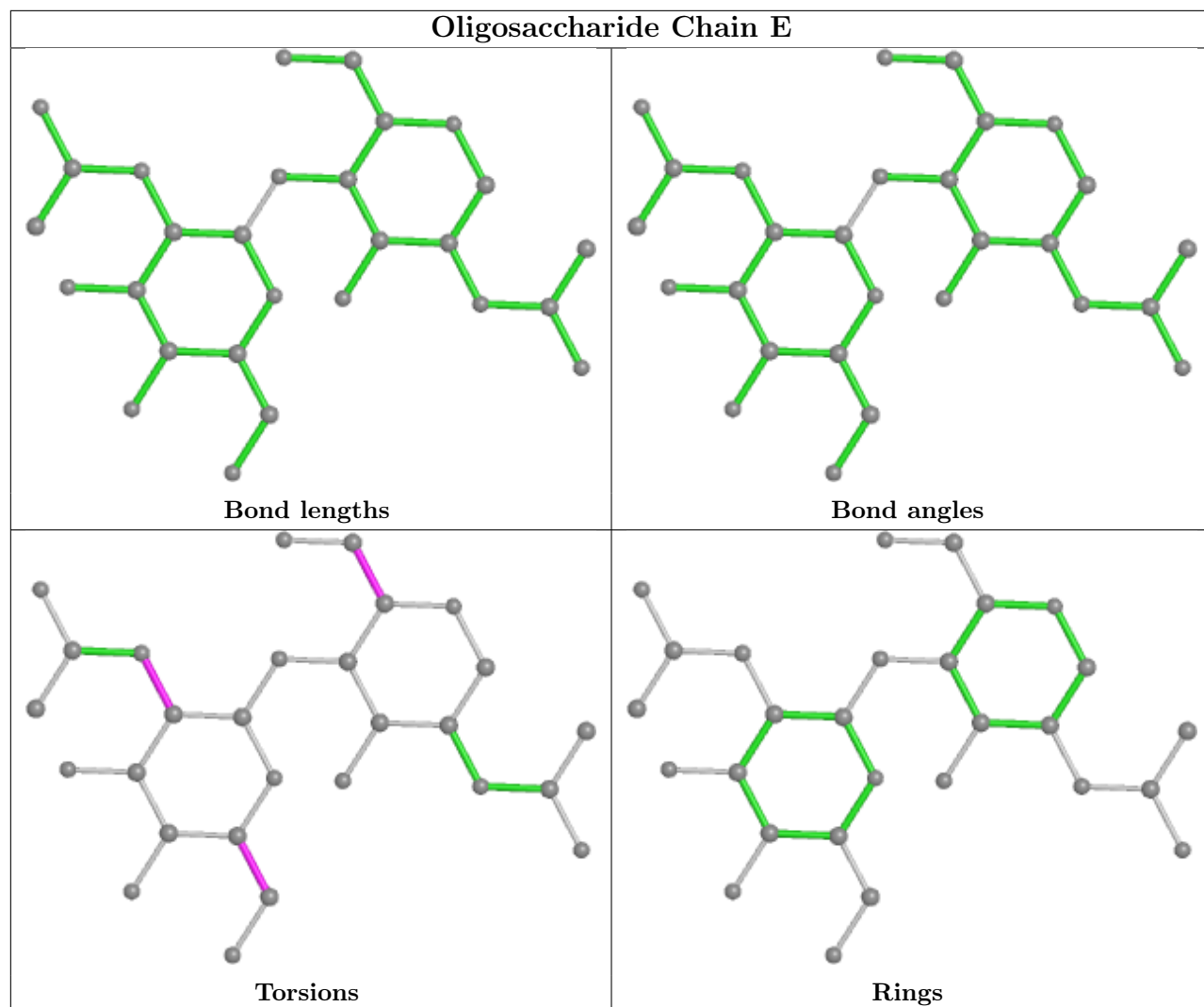
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	2	NAG	1	0
8	K	1	NAG	2	0
5	M	1	NAG	2	0
7	D	1	NAG	1	0
6	C	1	NAG	1	0

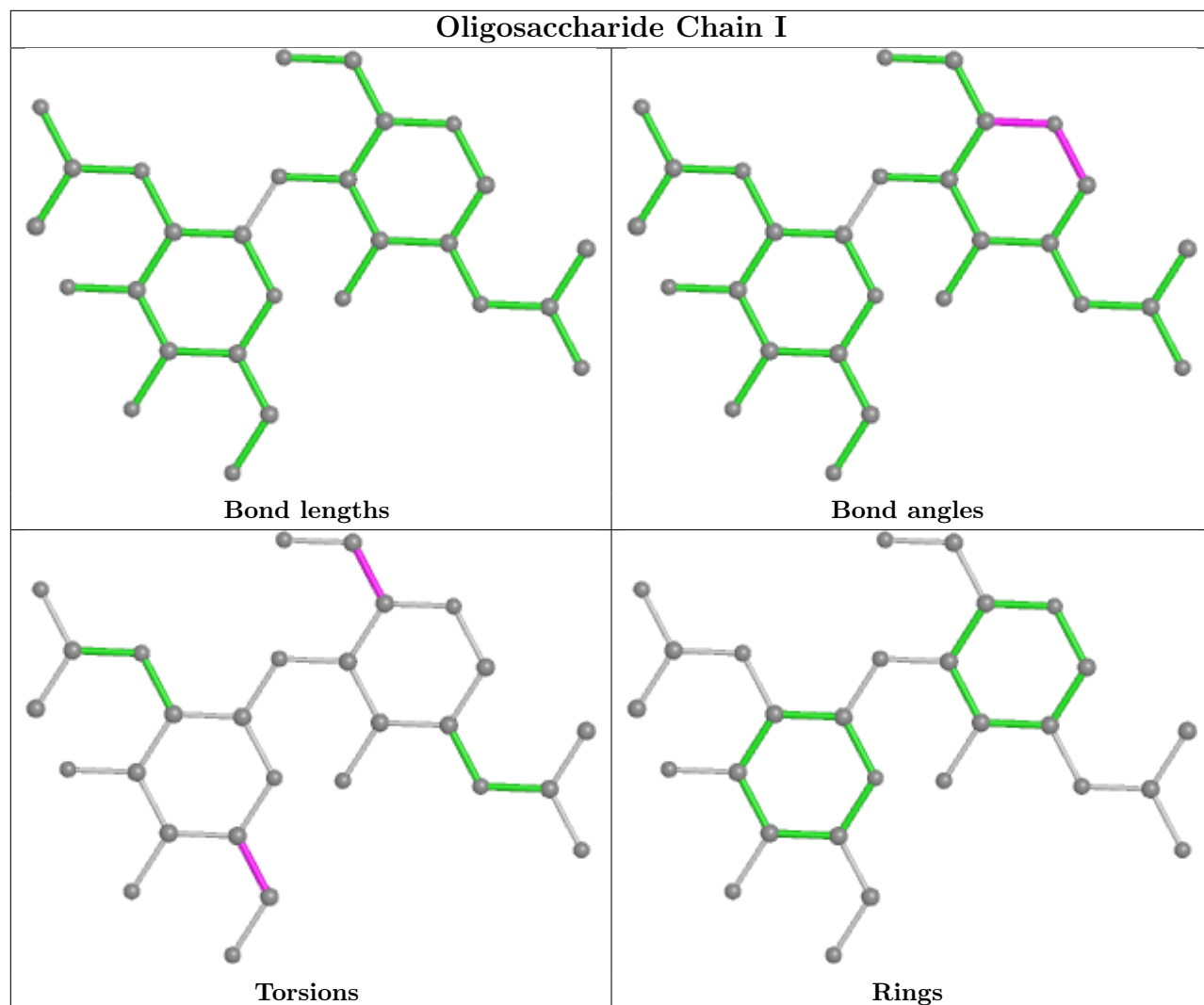
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



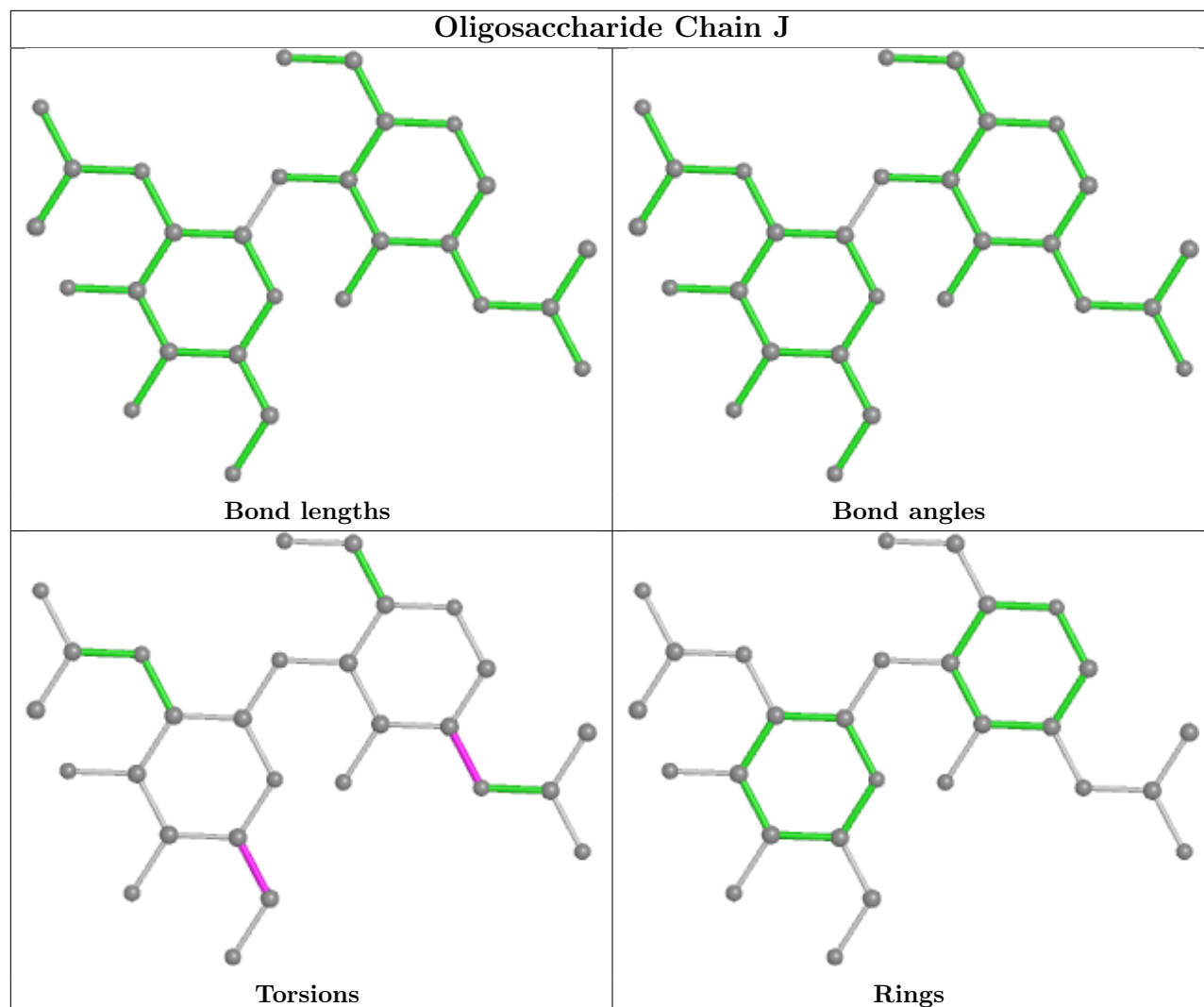


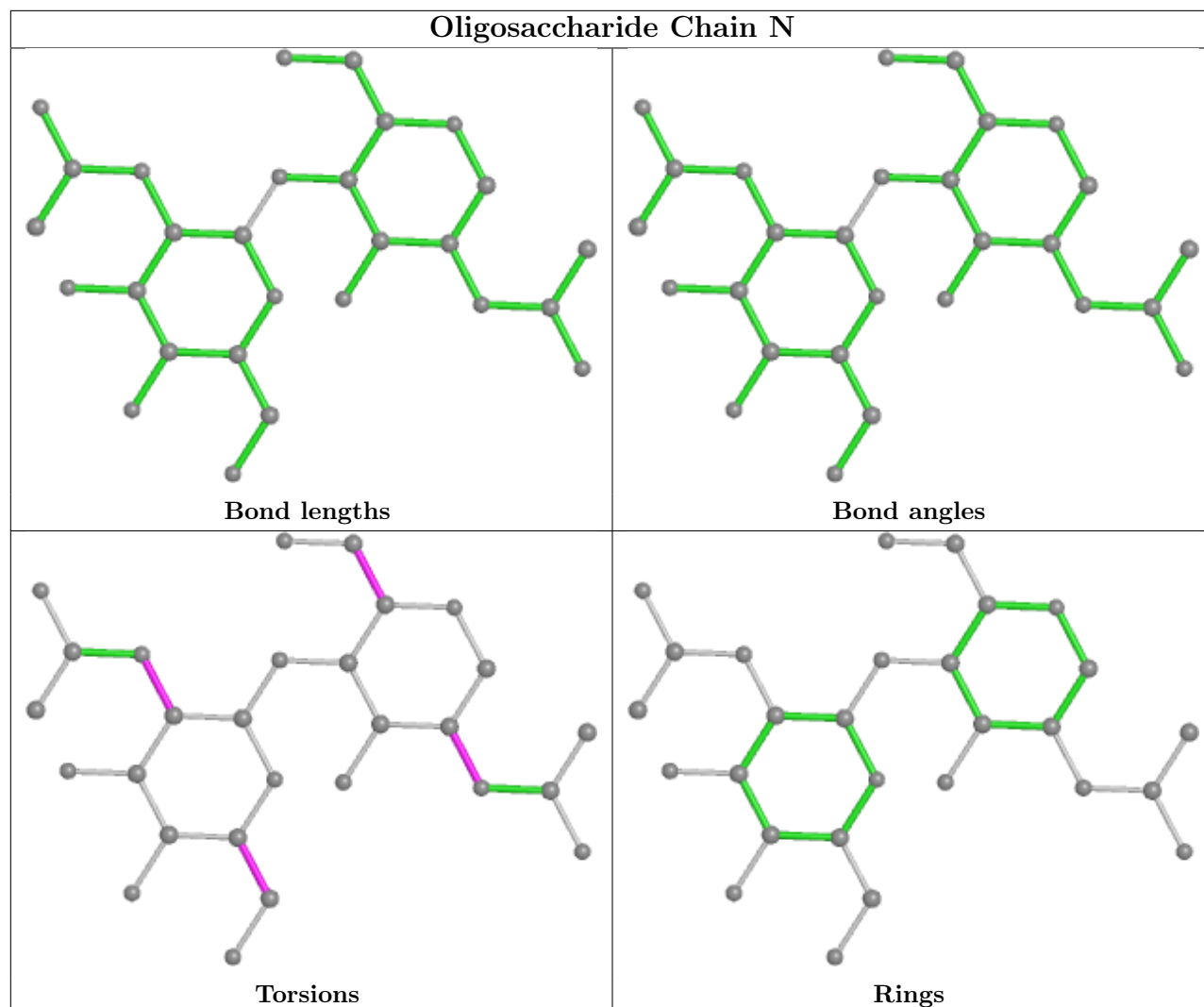


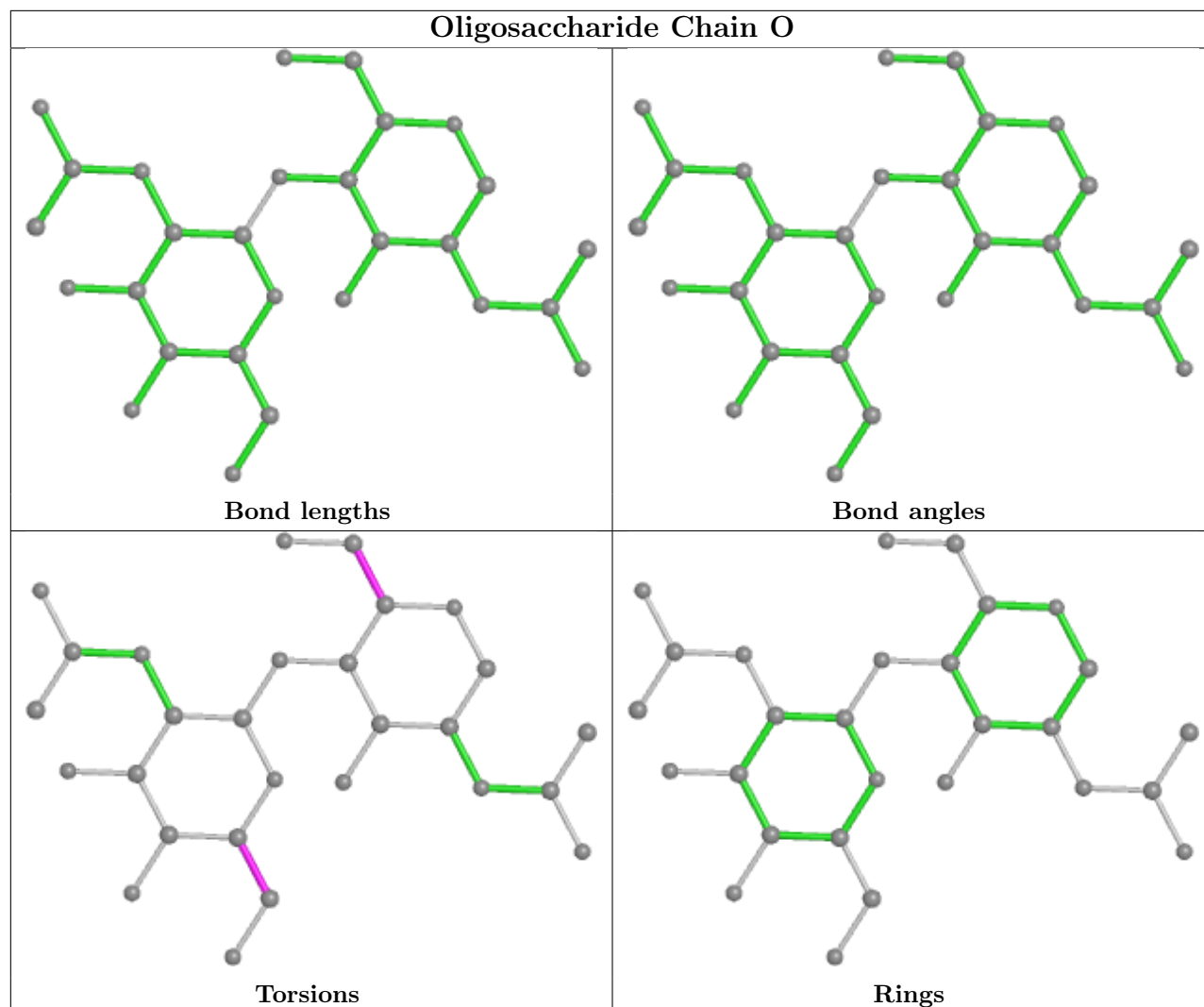


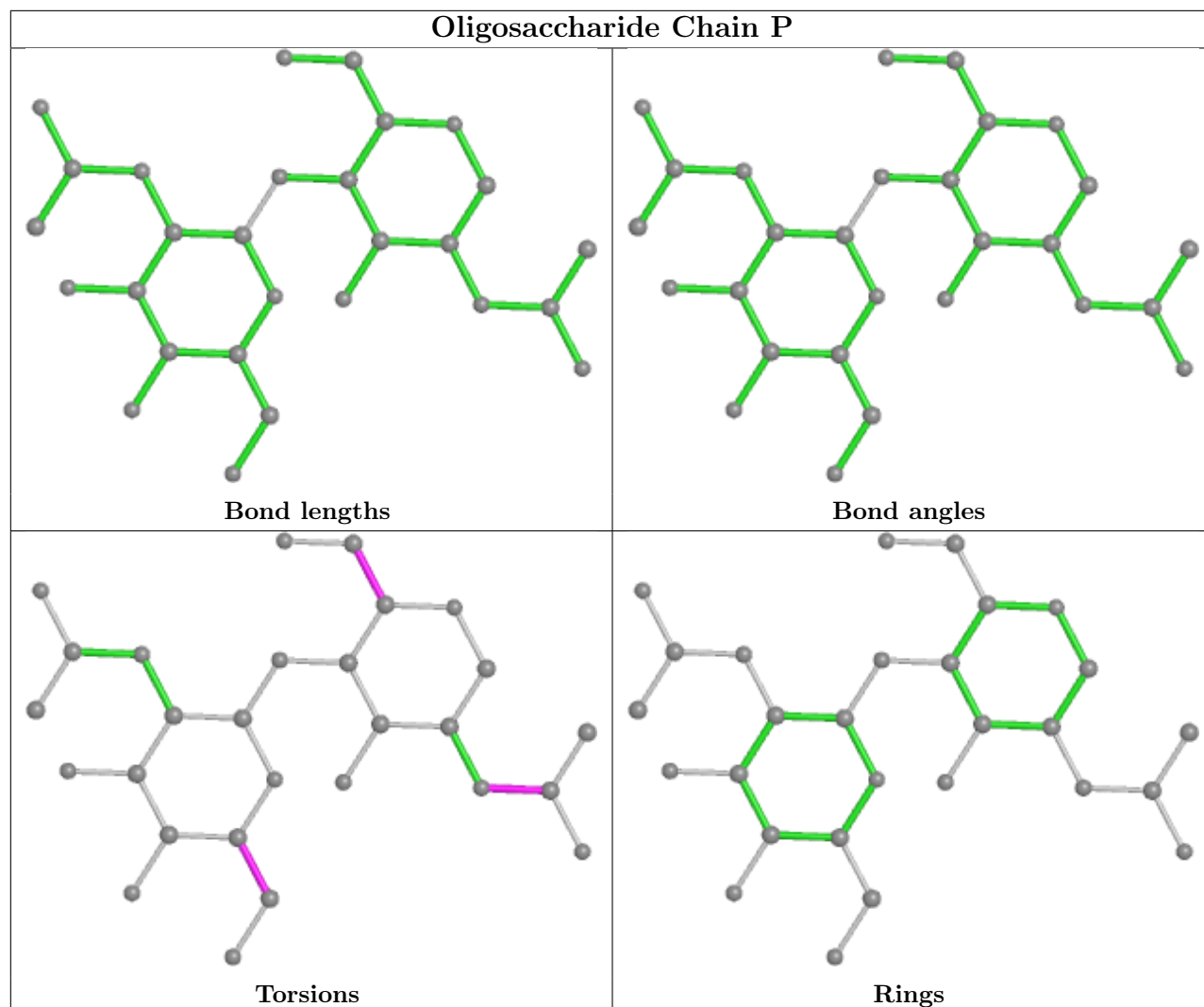


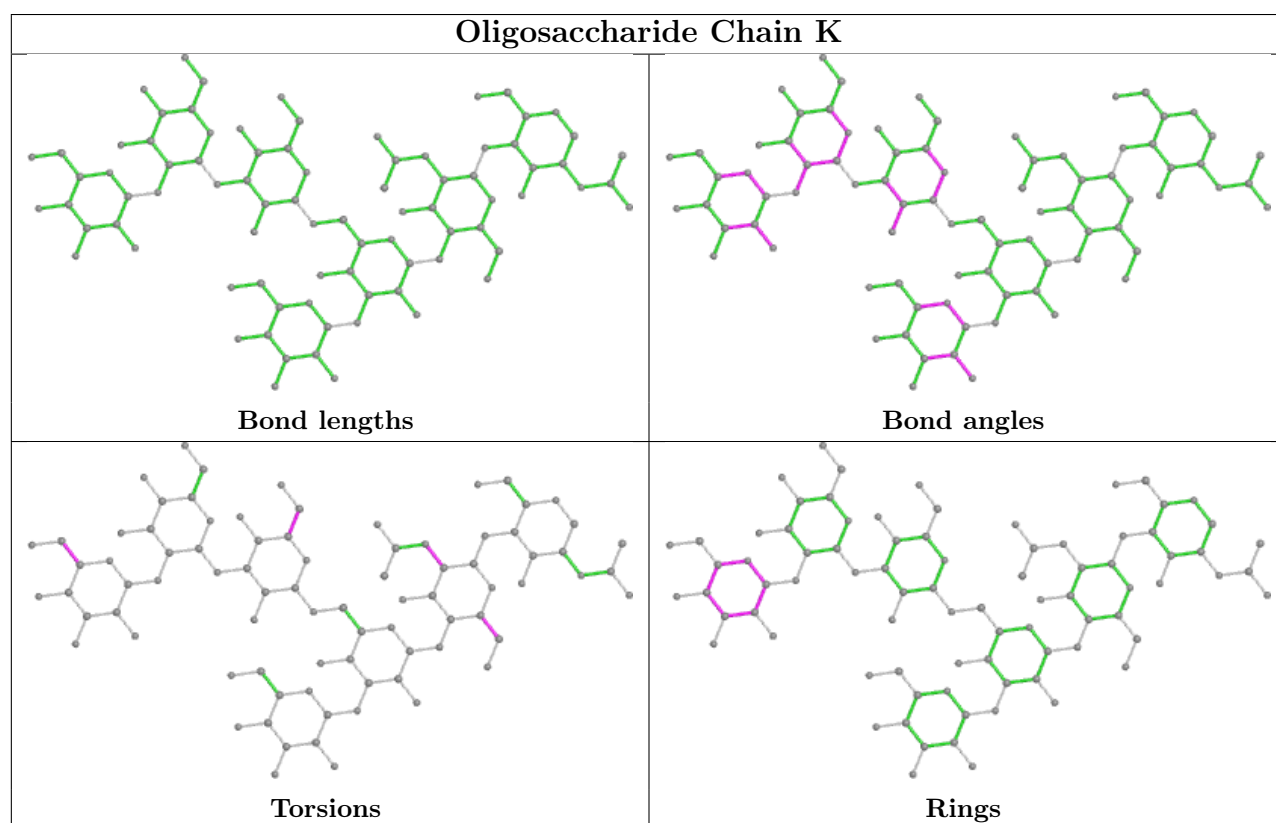
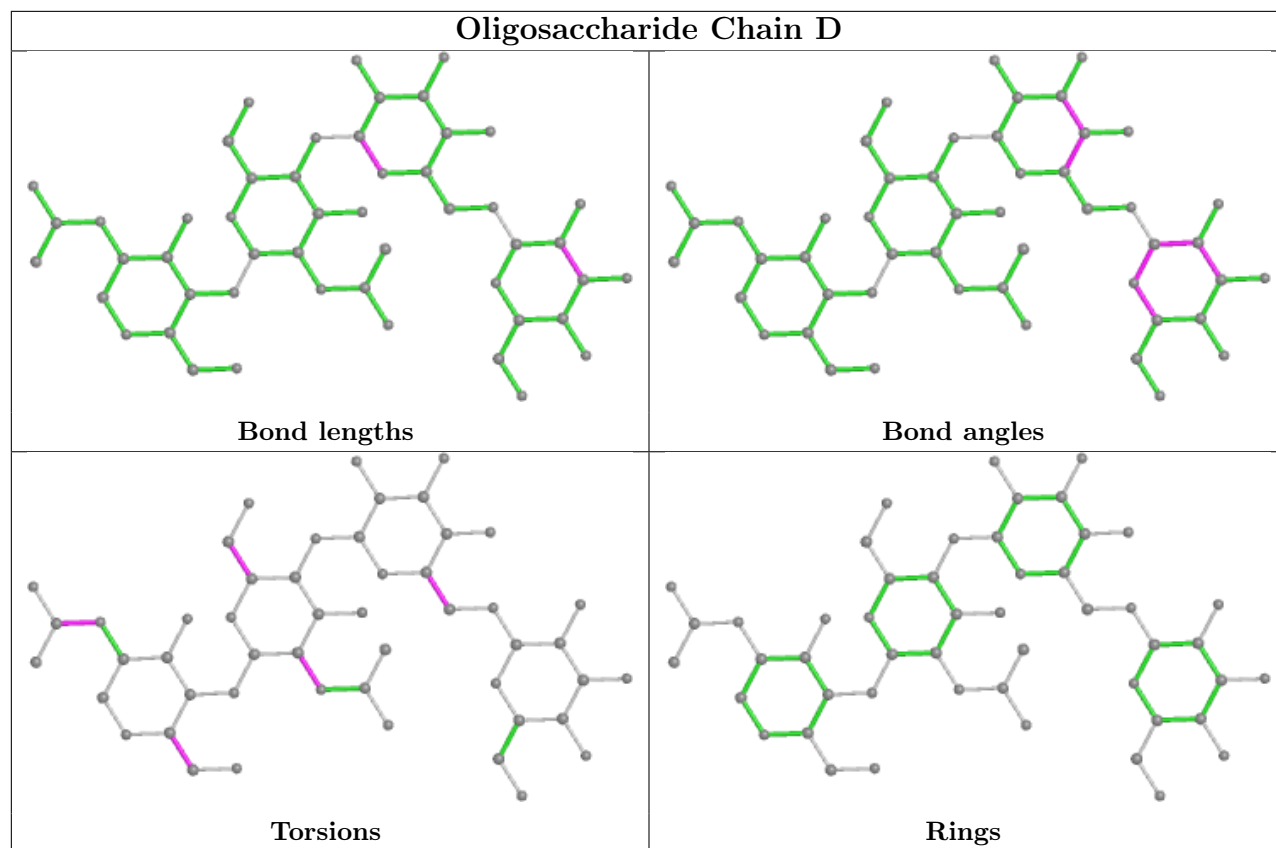












## 5.6 Ligand geometry

7 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
9	NAG	G	603	1	14,14,15	0.21	0	17,19,21	0.53	0
9	NAG	B	702	2	14,14,15	0.38	0	17,19,21	0.42	0
9	NAG	G	601	1	14,14,15	0.40	0	17,19,21	0.40	0
9	NAG	G	602	1	14,14,15	0.51	0	17,19,21	1.39	2 (11%)
9	NAG	G	604	1	14,14,15	0.33	0	17,19,21	0.34	0
9	NAG	B	703	2	14,14,15	0.67	0	17,19,21	0.39	0
9	NAG	B	701	2	14,14,15	0.24	0	17,19,21	0.33	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	NAG	G	603	1	-	4/6/23/26	0/1/1/1
9	NAG	B	702	2	-	0/6/23/26	0/1/1/1
9	NAG	G	601	1	-	2/6/23/26	0/1/1/1
9	NAG	G	602	1	-	5/6/23/26	0/1/1/1
9	NAG	G	604	1	-	2/6/23/26	0/1/1/1
9	NAG	B	703	2	-	0/6/23/26	0/1/1/1
9	NAG	B	701	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	G	602	NAG	C2-N2-C7	4.67	129.56	122.90
9	G	602	NAG	C1-C2-N2	2.23	114.30	110.49

There are no chirality outliers.

5 of 15 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	G	601	NAG	C4-C5-C6-O6
9	G	601	NAG	O5-C5-C6-O6
9	G	603	NAG	O5-C5-C6-O6
9	G	604	NAG	O5-C5-C6-O6
9	B	701	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B	702	NAG	1	0
9	G	602	NAG	4	0
9	B	703	NAG	1	0

## 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	G	449/481 (93%)	-0.06	10 (2%) 62 52	154, 213, 266, 325	0
2	B	122/153 (79%)	0.07	8 (6%) 18 14	215, 289, 343, 400	0
3	H	225/229 (98%)	-0.12	5 (2%) 62 52	149, 190, 243, 257	0
4	L	200/205 (97%)	0.03	10 (5%) 28 24	168, 212, 245, 265	0
All	All	996/1068 (93%)	-0.04	33 (3%) 46 37	149, 213, 298, 400	0

The worst 5 of 33 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	496	VAL	5.1
1	G	498	PRO	5.0
1	G	497	ALA	4.8
2	B	541	ALA	4.5
2	B	544	LEU	4.4

### 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, 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
5	BMA	A	3	11/12	0.63	0.52	336,340,341,342	0
8	MAN	K	4	11/12	0.65	0.24	389,393,397,398	0

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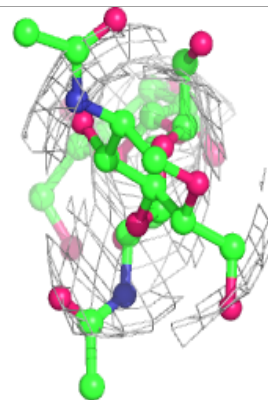
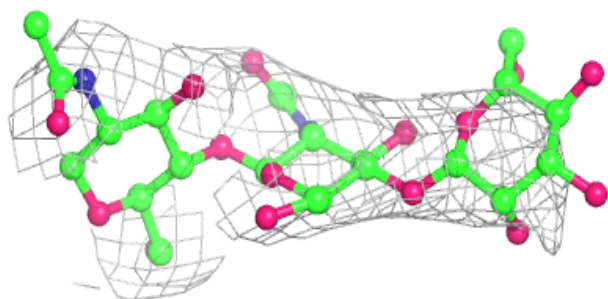
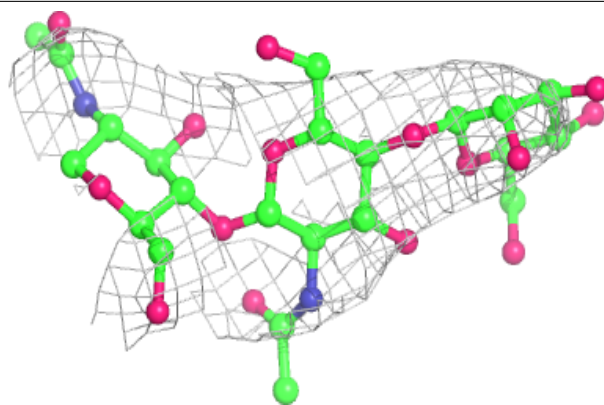
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	BMA	M	3	11/12	0.66	0.40	327,331,336,342	0
8	MAN	K	7	11/12	0.69	0.36	372,380,398,404	0
8	BMA	K	3	11/12	0.70	0.19	371,384,392,397	0
6	NAG	O	2	14/15	0.72	0.33	344,351,355,359	0
6	NAG	I	1	14/15	0.72	0.50	304,313,328,340	0
7	MAN	D	4	11/12	0.75	0.75	321,323,326,327	0
6	NAG	I	2	14/15	0.77	0.27	350,360,365,365	0
7	BMA	D	3	11/12	0.78	0.48	305,308,316,318	0
5	BMA	F	3	11/12	0.79	0.28	333,336,338,338	0
6	NAG	C	2	14/15	0.79	0.41	302,310,313,314	0
6	NAG	N	2	14/15	0.79	0.16	323,332,342,345	0
6	NAG	E	2	14/15	0.79	1.34	357,364,370,372	0
6	NAG	P	2	14/15	0.79	0.27	390,405,416,432	0
6	NAG	J	2	14/15	0.82	0.32	336,346,350,351	0
5	NAG	M	2	14/15	0.83	0.23	322,327,334,338	0
6	NAG	O	1	14/15	0.84	0.35	300,315,324,335	0
6	NAG	C	1	14/15	0.85	0.38	267,275,284,298	0
8	MAN	K	6	11/12	0.86	0.15	378,386,389,389	0
8	NAG	K	1	14/15	0.86	0.68	298,311,319,334	0
8	MAN	K	5	11/12	0.88	0.21	390,397,401,402	0
5	NAG	A	2	14/15	0.88	0.22	306,314,321,328	0
6	NAG	P	1	14/15	0.88	0.33	321,332,353,373	0
8	NAG	K	2	14/15	0.91	0.35	340,348,353,363	0
7	NAG	D	2	14/15	0.91	0.29	291,296,304,307	0
6	NAG	E	1	14/15	0.91	0.50	312,326,341,347	0
5	NAG	A	1	14/15	0.92	0.20	270,282,289,296	0
5	NAG	M	1	14/15	0.93	0.26	279,288,299,310	0
6	NAG	J	1	14/15	0.93	0.20	293,308,315,329	0
7	NAG	D	1	14/15	0.94	0.24	253,270,277,284	0
5	NAG	F	2	14/15	0.95	0.22	318,327,333,334	0
5	NAG	F	1	14/15	0.96	0.32	281,290,303,311	0
6	NAG	N	1	14/15	0.96	0.22	279,291,313,318	0

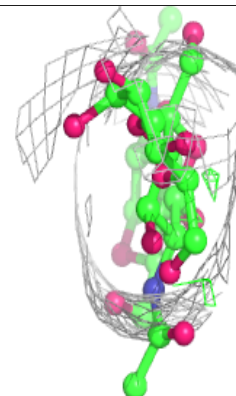
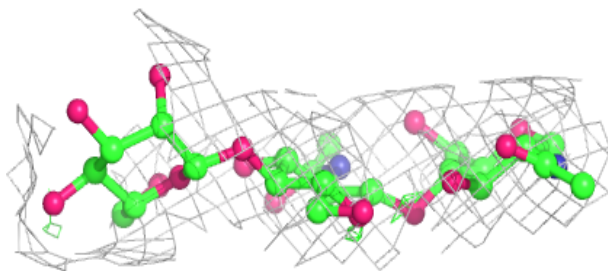
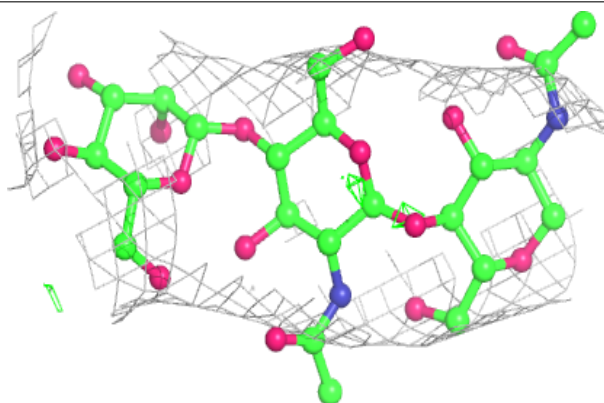
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

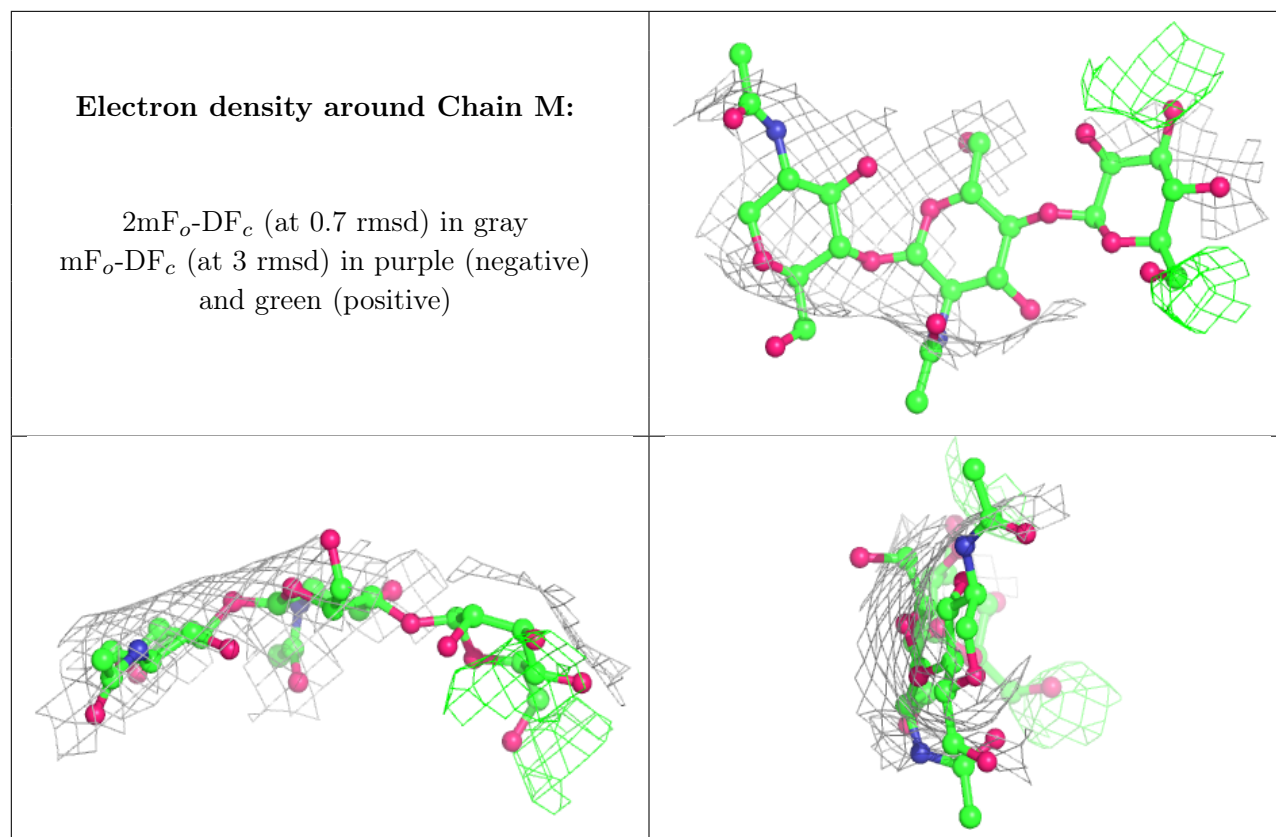
**Electron density around Chain A:**

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

**Electron density around Chain F:**

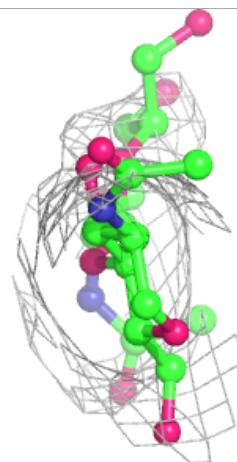
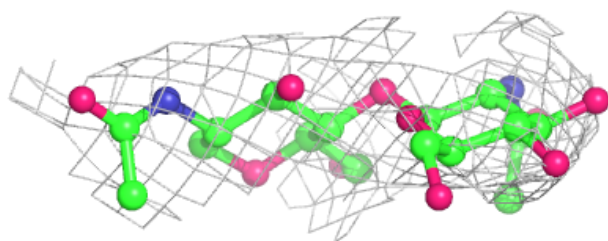
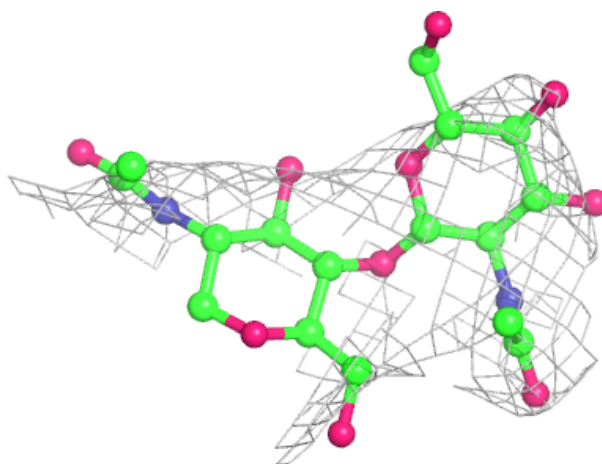
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





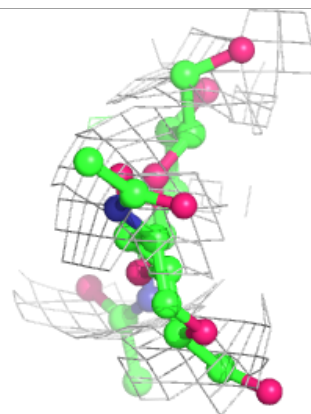
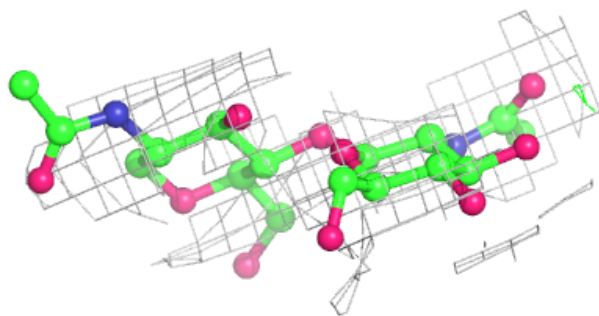
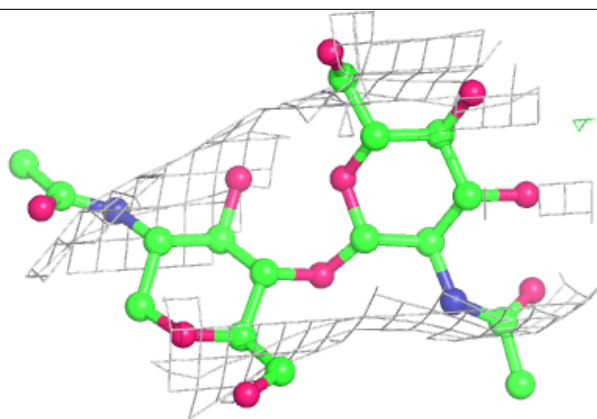
**Electron density around Chain C:**

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

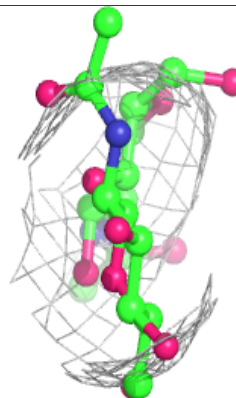
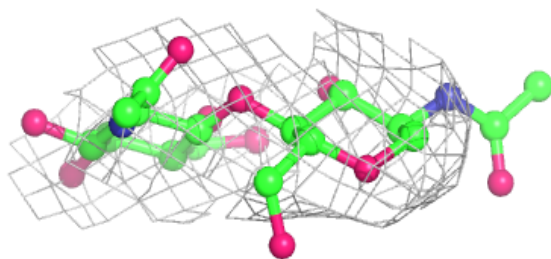
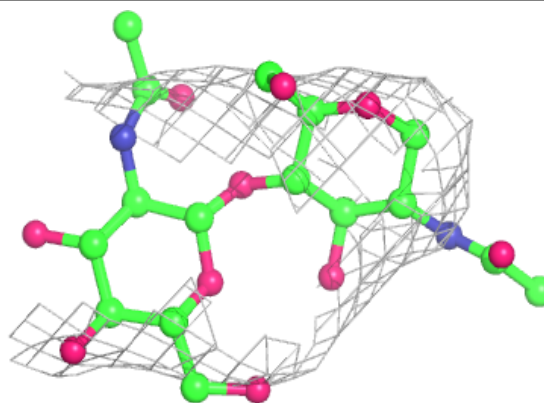


**Electron density around Chain E:**

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

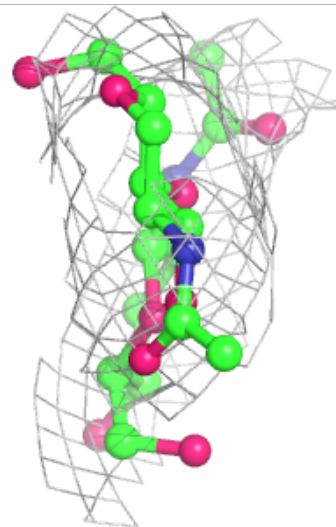
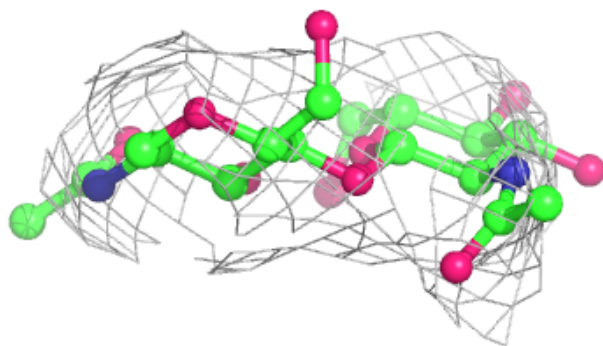
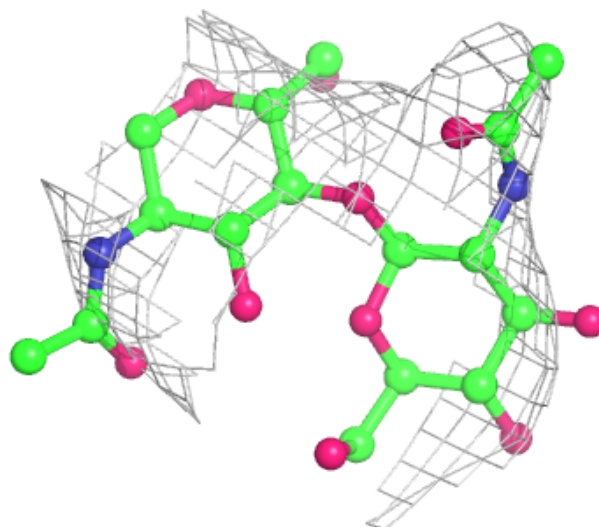
**Electron density around Chain I:**

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



**Electron density around Chain J:**

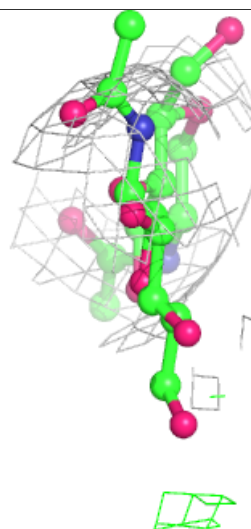
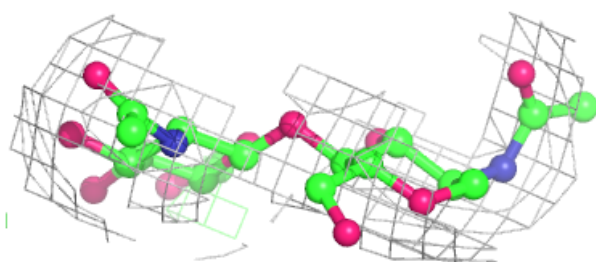
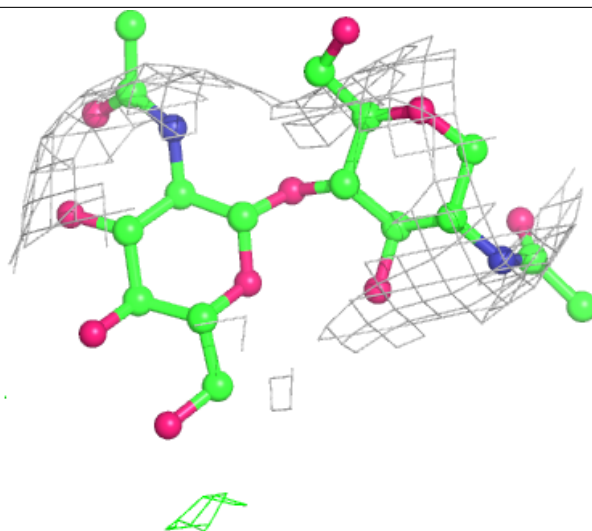
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





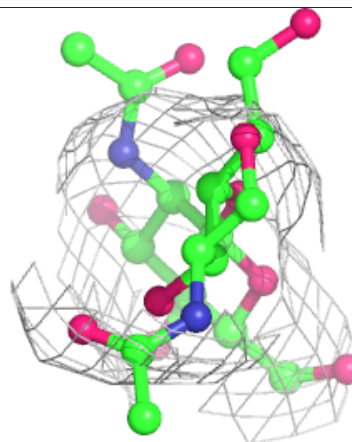
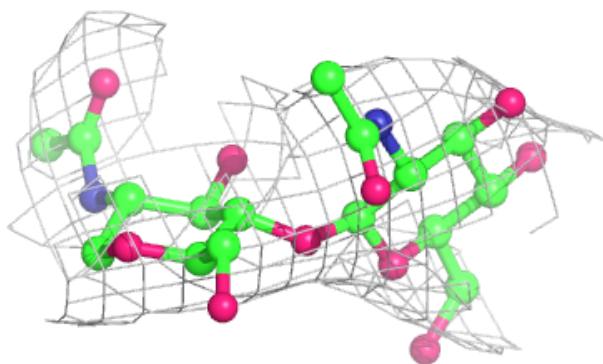
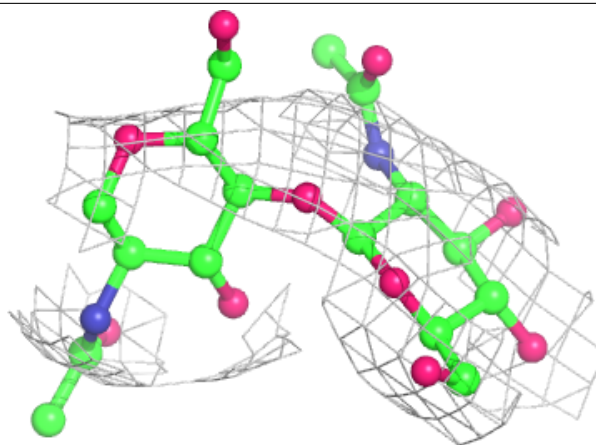
**Electron density around Chain N:**

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

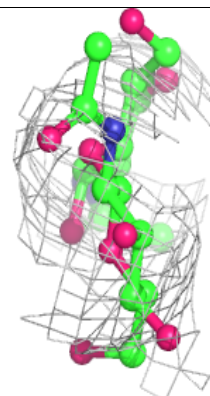
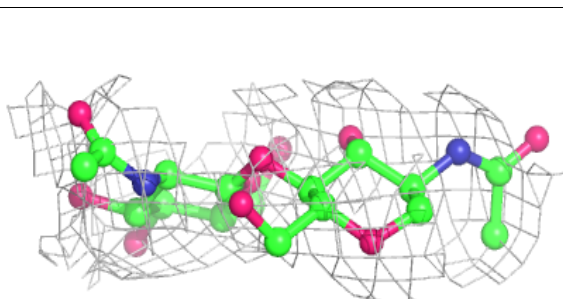
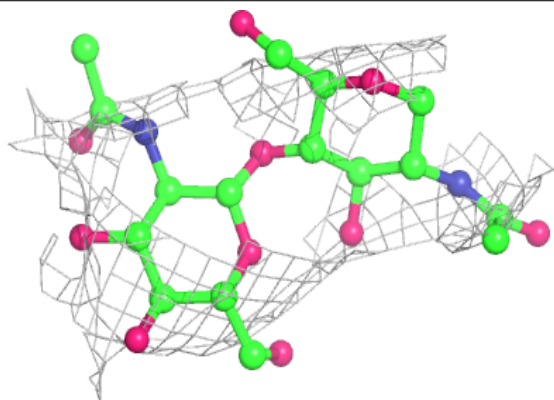


**Electron density around Chain O:**

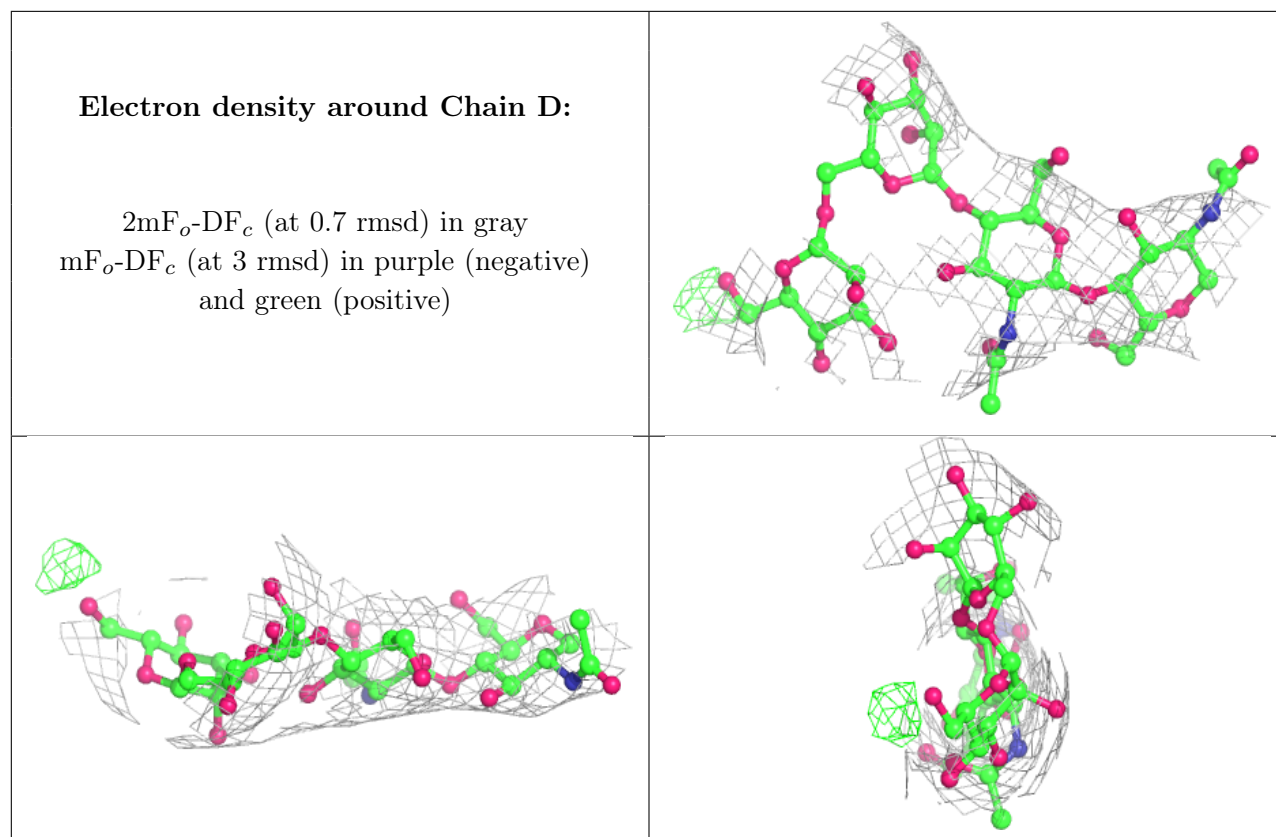
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

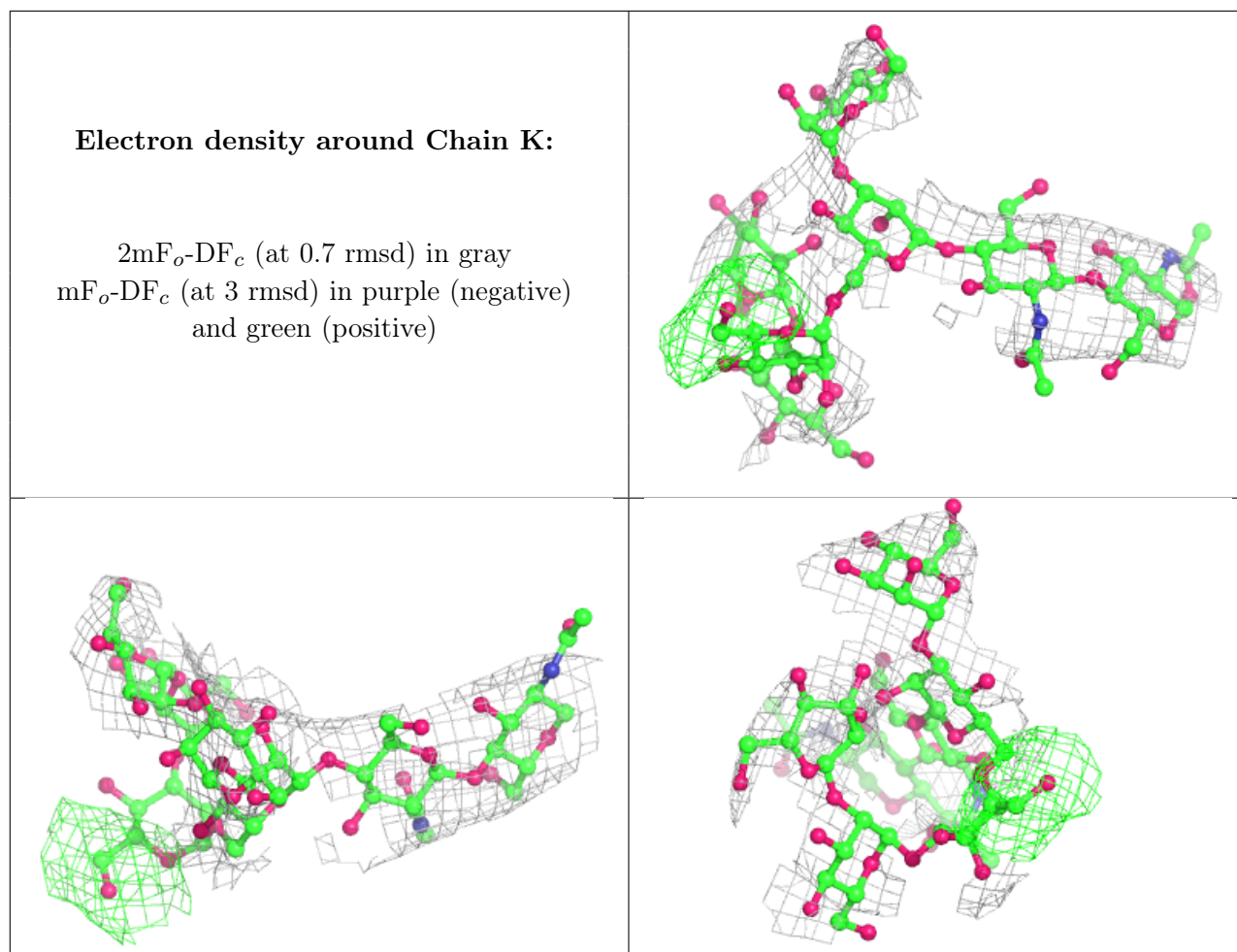
**Electron density around Chain P:**

$2mF_o-DF_c$  (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
9	NAG	G	603	14/15	0.80	0.53	341,352,358,360	0
9	NAG	B	703	14/15	0.82	0.18	200,230,258,262	0
9	NAG	B	702	14/15	0.83	0.20	221,236,260,261	0
9	NAG	G	604	14/15	0.83	0.26	321,334,340,341	0
9	NAG	G	601	14/15	0.85	0.30	262,271,275,276	0
9	NAG	G	602	14/15	0.86	0.36	293,305,313,314	0
9	NAG	B	701	14/15	0.86	0.47	212,234,263,273	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.