



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 8, 2020 – 02:18 PM BST

PDB ID : 4UO0
Title : Structure of the A_Equine_Richmond_07 H3 haemagglutinin
Authors : Vachieri, S.G.; Collins, P.J.; Haire, L.F.; Ogradowicz, R.W.; Martin, S.R.;
Walker, P.A.; Xiong, X.; Gamblin, S.J.; Skehel, J.J.
Deposited on : 2014-05-31
Resolution : 1.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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 : 2.13.1
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) : 2.13.1

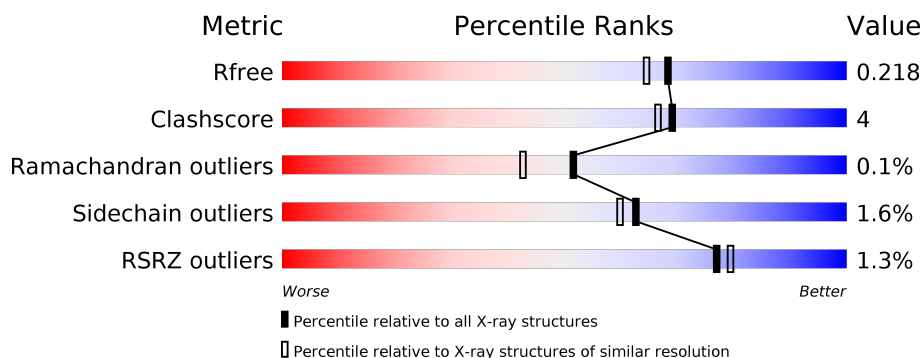
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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.



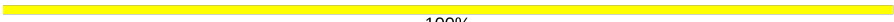

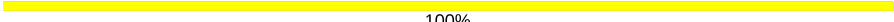



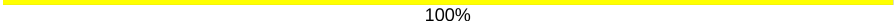
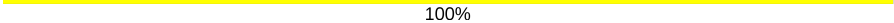
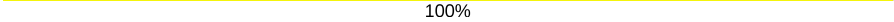



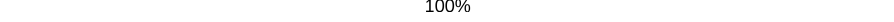

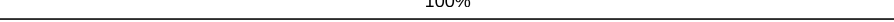


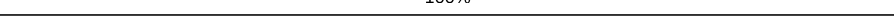
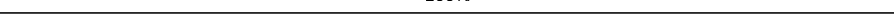



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	329	<div> <div>2%</div> <div> <div></div> <div>91%</div> <div>7%</div> <div>•</div> </div> </div>
1	C	329	<div> <div>%</div> <div> <div></div> <div>92%</div> <div>5%</div> <div>••</div> </div> </div>
1	E	329	<div> <div>%</div> <div> <div></div> <div>91%</div> <div>7%</div> <div>••</div> </div> </div>
2	B	172	<div> <div>%</div> <div> <div></div> <div>92%</div> <div>8%</div> </div> </div>
2	D	172	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div>•</div> </div> </div>
2	F	172	<div> <div>%</div> <div> <div></div> <div>92%</div> <div>7%</div> <div>•</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	G	2	 100%
3	H	2	 50% 50%
3	J	2	 100%
3	L	2	 50% 50%
3	R	2	 100%
3	V	2	 100%
3	W	2	 100%
3	Y	2	 100%
3	a	2	 100%
4	I	5	 60% 40%
4	P	5	 40% 60%
4	Q	5	 20% 80%
4	X	5	 100%
5	K	4	 75% 25%
6	M	3	 100%
6	O	3	 33% 67%
6	T	3	 67% 33%
7	N	2	 100%
8	S	6	 100%
9	U	3	 100%
9	Z	3	 33% 67%
9	b	3	 33% 67%

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
10	BMA	A	604	-	-	-	X
12	NAG	C	601	-	-	-	X
12	NAG	E	621	-	-	-	X
3	NAG	G	2	-	-	-	X
3	NAG	H	2	-	-	-	X
3	NAG	L	1	X	-	-	-
3	NAG	W	2	-	-	-	X
3	NAG	Y	2	-	-	-	X
4	MAN	P	4	-	-	-	X
4	MAN	P	5	-	-	-	X
5	FUC	K	4	X	-	-	-
7	FUC	N	2	X	-	-	-
8	FUC	S	6	X	-	-	-
9	FUC	U	3	X	-	-	-
9	FUC	Z	3	X	-	-	-
9	FUC	b	3	X	-	-	-

2 Entry composition [i](#)

There are 13 unique types of molecules in this entry. The entry contains 14213 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 HEMAGGLUTININ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	8	0
			2573	1612	454	492	15			
1	C	320	Total	C	N	O	S	0	1	0
			2496	1560	441	480	15			
1	E	325	Total	C	N	O	S	0	4	0
			2548	1593	449	491	15			

- Molecule 2 is a protein called HEMAGGLUTININ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	172	Total	C	N	O	S	0	6	0
			1428	896	242	283	7			
2	D	172	Total	C	N	O	S	0	4	0
			1421	890	245	279	7			
2	F	172	Total	C	N	O	S	0	7	0
			1437	903	245	281	8			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



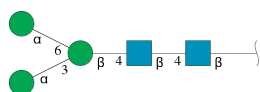
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	J	2	Total	C	N	O	0	0	0
			28	16	2	10			

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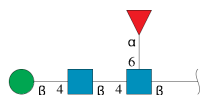
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	L	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	R	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	V	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	W	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	Y	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	a	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



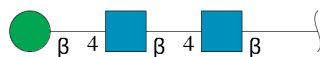
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	I	5	Total	C	N	O	0	0	0
			61	34	2	25			
4	P	5	Total	C	N	O	0	0	0
			61	34	2	25			
4	Q	5	Total	C	N	O	0	0	0
			61	34	2	25			
4	X	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	K	4	Total	C	N	O	0	0	0
			49	28	2	19			

- Molecule 6 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.



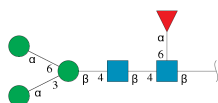
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	M	3	Total	C	N	O	0	0	0
			39	22	2	15			
6	O	3	Total	C	N	O	0	0	0
			39	22	2	15			
6	T	3	Total	C	N	O	0	0	0
			39	22	2	15			

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



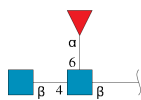
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	N	2	Total	C	N	O	0	0	0
			24	14	1	9			

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



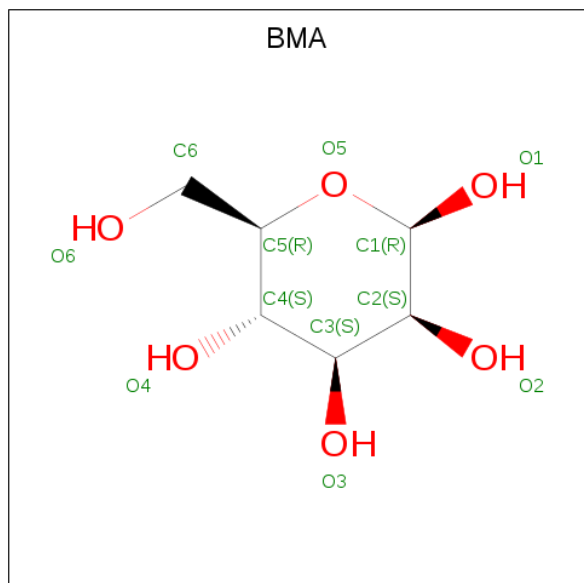
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
8	S	6	Total	C	N	O	0	0	0
			71	40	2	29			

- Molecule 9 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	U	3	Total	C	N	O	0	0	0
			38	22	2	14			
9	Z	3	Total	C	N	O	0	0	0
			38	22	2	14			
9	b	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 10 is beta-D-mannopyranose (three-letter code: BMA) (formula: C₆H₁₂O₆).



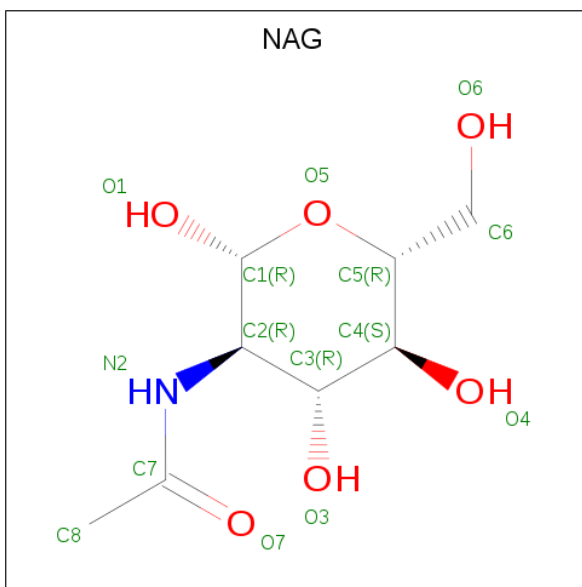
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			11	6	5		

- Molecule 11 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	A	1	Total	C	O	0	0
			4	2	2		
11	A	1	Total	C	O	0	0
			4	2	2		
11	B	1	Total	C	O	0	0
			4	2	2		
11	B	1	Total	C	O	0	0
			4	2	2		
11	C	1	Total	C	O	0	0
			4	2	2		
11	C	1	Total	C	O	0	0
			4	2	2		
11	D	1	Total	C	O	0	0
			4	2	2		
11	E	1	Total	C	O	0	0
			4	2	2		
11	E	1	Total	C	O	0	0
			4	2	2		
11	F	1	Total	C	O	0	0
			4	2	2		
11	F	1	Total	C	O	0	0
			4	2	2		
11	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 12 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	C	1	Total	C	N	O	0	0
			14	8	1	5		
12	E	1	Total	C	N	O	0	0
			14	8	1	5		

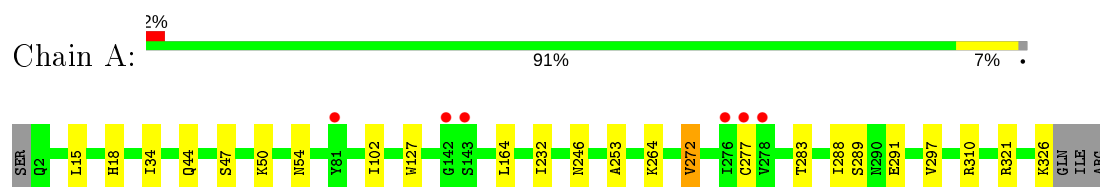
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	314	Total	O	0	0
			314	314		
13	B	191	Total	O	0	0
			191	191		
13	C	283	Total	O	0	0
			283	283		
13	D	174	Total	O	0	0
			174	174		
13	E	256	Total	O	0	0
			256	256		
13	F	134	Total	O	0	0
			134	134		

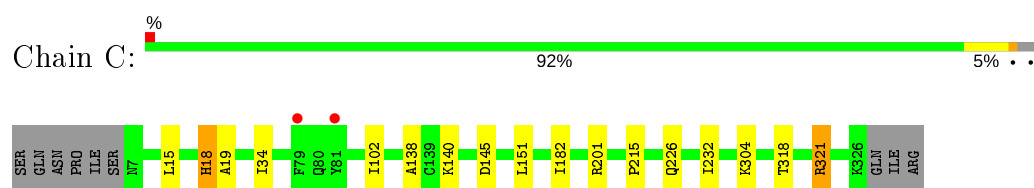
3 Residue-property plots [i](#)

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.

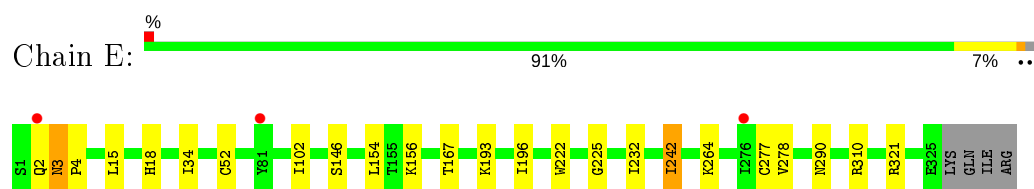
- Molecule 1: HEMAGGLUTININ



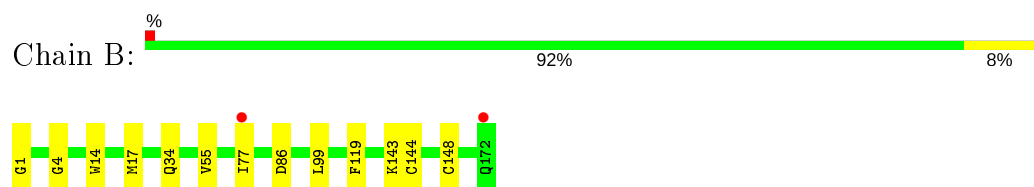
- Molecule 1: HEMAGGLUTININ



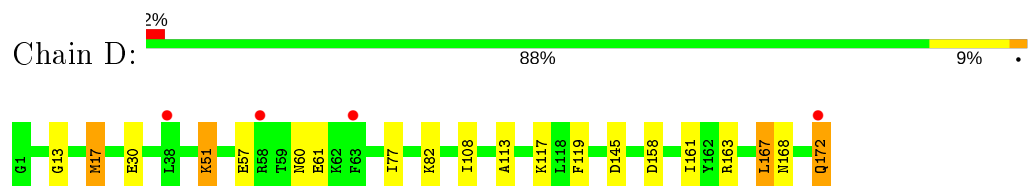
- Molecule 1: HEMAGGLUTININ



- Molecule 2: HEMAGGLUTININ



- Molecule 2: HEMAGGLUTININ



- Molecule 2: HEMAGGLUTININ



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1
MAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain W:  100%MAG1
MAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Y:  100%MAG1
MAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain a:  100%MAG1
MAG2


- Molecule 4: alpha-D-mannopyranose-(1-3)-[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 I:  60% 40%MAG1
MAG2
MAN3
MAN4
MAN5

- Molecule 4: alpha-D-mannopyranose-(1-3)-[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

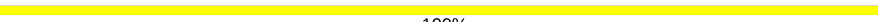
Chain P:  40% 60%MAG1
MAG2
MAN3
MAN4
MAN5

- Molecule 4: alpha-D-mannopyranose-(1-3)-[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 Q:  20% 80%




- Molecule 4: alpha-D-mannopyranose-(1-3)-[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 X:  100%



- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  75% 25%



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

Chain M:  100%



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

Chain O:  33% 67%



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

Chain T:  67% 33%



- Molecule 7: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  100%



- Molecule 8: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-[α -L-fucopyranose-(1-6)]2-acetamido-2-deoxy- β -D-glucopyranose

Chain S:  100%

MAG1
MAG2
BMA3
MAN4
MAN5
FUC6

- Molecule 9: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-[α -L-fucopyranose-(1-6)]2-acetamido-2-deoxy- β -D-glucopyranose

Chain U:  100%

MAG1
MAG2
FUC3

- Molecule 9: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-[α -L-fucopyranose-(1-6)]2-acetamido-2-deoxy- β -D-glucopyranose

Chain Z:  33%  67%

MAG1
MAG2
FUC3

- Molecule 9: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-[α -L-fucopyranose-(1-6)]2-acetamido-2-deoxy- β -D-glucopyranose

Chain b:  33%  67%

MAG1
MAG2
FUC3

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.41Å 129.33Å 192.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	107.37 – 1.90 46.58 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.3 (107.37-1.90) 97.3 (46.58-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.191 , 0.218 0.191 , 0.218	Depositor DCC
R_{free} test set	7614 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.661	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14213	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FUC, MAN, BMA, NAG, EDO

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	0.39	0/2651	0.56	0/3598
1	C	0.41	0/2552	0.58	0/3464
1	E	0.39	0/2614	0.57	0/3551
2	B	0.44	0/1471	0.57	0/1978
2	D	0.42	0/1458	0.59	0/1960
2	F	0.44	0/1483	0.61	0/1994
All	All	0.41	0/12229	0.58	0/16545

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	A	2573	0	2547	16	0
1	C	2496	0	2440	14	0
1	E	2548	0	2499	16	0
2	B	1428	0	1363	25	0
2	D	1421	0	1360	19	0
2	F	1437	0	1386	19	0
3	G	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	28	0	25	1	0
3	J	28	0	25	0	0
3	L	28	0	25	0	0
3	R	28	0	25	0	0
3	V	28	0	25	1	0
3	W	28	0	25	0	0
3	Y	28	0	25	0	0
3	a	28	0	25	0	0
4	I	61	0	52	2	0
4	P	61	0	52	4	0
4	Q	61	0	52	1	0
4	X	61	0	52	0	0
5	K	49	0	43	2	0
6	M	39	0	34	0	0
6	O	39	0	34	0	0
6	T	39	0	34	1	0
7	N	24	0	22	0	0
8	S	71	0	61	0	0
9	U	38	0	34	0	0
9	Z	38	0	34	0	0
9	b	38	0	34	0	0
10	A	11	0	10	0	0
11	A	8	0	12	1	0
11	B	8	0	12	0	0
11	C	8	0	12	0	0
11	D	4	0	6	0	0
11	E	8	0	12	1	0
11	F	12	0	18	0	0
12	C	14	0	13	0	0
12	E	14	0	13	0	0
13	A	314	0	0	4	0
13	B	191	0	0	2	0
13	C	283	0	0	0	0
13	D	174	0	0	2	0
13	E	256	0	0	1	0
13	F	134	0	0	2	0
All	All	14213	0	12466	93	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 4.

The worst 5 of 93 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)
2:B:77[A]:ILE:HD12	2:F:77[A]:ILE:CD1	2.03	0.88
2:B:77[A]:ILE:HD12	2:F:77[A]:ILE:HD11	1.58	0.83
2:B:77[A]:ILE:CD1	2:F:77[A]:ILE:CD1	2.57	0.82
1:E:167:THR:OG1	1:E:242:ILE:HD11	1.80	0.81
4:I:3:BMA:H62	4:I:5:MAN:H5	1.68	0.76

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	331/329 (101%)	323 (98%)	8 (2%)	0	100	100
1	C	319/329 (97%)	309 (97%)	10 (3%)	0	100	100
1	E	327/329 (99%)	314 (96%)	10 (3%)	3 (1%)	17	7
2	B	176/172 (102%)	167 (95%)	9 (5%)	0	100	100
2	D	174/172 (101%)	166 (95%)	8 (5%)	0	100	100
2	F	177/172 (103%)	169 (96%)	8 (4%)	0	100	100
All	All	1504/1503 (100%)	1448 (96%)	53 (4%)	3 (0%)	51	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	2	GLN
1	E	3[A]	ASN
1	E	3[B]	ASN

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	296/292 (101%)	292 (99%)	4 (1%)	67	65
1	C	283/292 (97%)	280 (99%)	3 (1%)	73	73
1	E	291/292 (100%)	287 (99%)	4 (1%)	67	65
2	B	151/146 (103%)	151 (100%)	0	100	100
2	D	150/146 (103%)	142 (95%)	8 (5%)	22	13
2	F	153/146 (105%)	149 (97%)	4 (3%)	46	39
All	All	1324/1314 (101%)	1301 (98%)	23 (2%)	62	57

5 of 23 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	57	GLU
2	D	108	ILE
2	F	57	GLU
2	D	60	ASN
2	D	167	LEU

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

Mol	Chain	Res	Type
1	A	54	ASN
1	A	188	ASN
1	C	18	HIS
1	C	296	ASN
2	F	172	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

68 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$
3	NAG	G	1	1,3	14,14,15	0.54	0	17,19,21	1.48	2 (11%)
3	NAG	G	2	3	14,14,15	0.53	0	17,19,21	1.03	1 (5%)
3	NAG	H	1	1,3	14,14,15	0.62	0	17,19,21	1.21	1 (5%)
3	NAG	H	2	3	14,14,15	0.55	0	17,19,21	0.84	0
4	NAG	I	1	1,4	14,14,15	0.66	0	17,19,21	1.05	1 (5%)
4	NAG	I	2	4	14,14,15	0.59	0	17,19,21	0.94	1 (5%)
4	BMA	I	3	4	11,11,12	0.63	0	15,15,17	2.75	5 (33%)
4	MAN	I	4	4	11,11,12	0.52	0	15,15,17	0.91	1 (6%)
4	MAN	I	5	4	11,11,12	0.60	0	15,15,17	1.40	2 (13%)
3	NAG	J	1	1,3	14,14,15	0.62	0	17,19,21	0.90	1 (5%)
3	NAG	J	2	3	14,14,15	0.67	0	17,19,21	1.66	3 (17%)
5	NAG	K	1	1,5	14,14,15	0.67	0	17,19,21	1.07	2 (11%)
5	NAG	K	2	5	14,14,15	0.59	0	17,19,21	1.25	3 (17%)
5	BMA	K	3	5	11,11,12	0.43	0	15,15,17	1.34	3 (20%)
5	FUC	K	4	5	10,10,11	0.65	0	14,14,16	1.00	1 (7%)
3	NAG	L	1	1,3	14,14,15	0.62	0	17,19,21	1.08	0
3	NAG	L	2	3	14,14,15	0.59	0	17,19,21	1.39	2 (11%)
6	NAG	M	1	1,6	14,14,15	0.46	0	17,19,21	1.10	2 (11%)
6	NAG	M	2	6	14,14,15	0.52	0	17,19,21	1.07	1 (5%)
6	BMA	M	3	6	11,11,12	0.42	0	15,15,17	0.94	1 (6%)
7	NAG	N	1	2,7	14,14,15	0.55	0	17,19,21	1.73	2 (11%)
7	FUC	N	2	7	10,10,11	0.65	0	14,14,16	1.06	2 (14%)
6	NAG	O	1	1,6	14,14,15	0.60	0	17,19,21	0.93	0
6	NAG	O	2	6	14,14,15	0.61	0	17,19,21	1.32	3 (17%)
6	BMA	O	3	6	11,11,12	0.67	0	15,15,17	2.41	4 (26%)
4	NAG	P	1	1,4	14,14,15	0.62	0	17,19,21	1.72	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	P	2	4	14,14,15	0.75	0	17,19,21	1.20	1 (5%)
4	BMA	P	3	4	11,11,12	0.50	0	15,15,17	0.98	1 (6%)
4	MAN	P	4	4	11,11,12	0.81	0	15,15,17	1.68	2 (13%)
4	MAN	P	5	4	11,11,12	0.52	0	15,15,17	1.59	2 (13%)
4	NAG	Q	1	1,4	14,14,15	0.60	0	17,19,21	0.90	0
4	NAG	Q	2	4	14,14,15	0.52	0	17,19,21	1.42	3 (17%)
4	BMA	Q	3	4	11,11,12	0.38	0	15,15,17	0.84	0
4	MAN	Q	4	4	11,11,12	0.59	0	15,15,17	0.90	1 (6%)
4	MAN	Q	5	4	11,11,12	0.61	0	15,15,17	1.29	2 (13%)
3	NAG	R	1	1,3	14,14,15	0.62	0	17,19,21	1.04	0
3	NAG	R	2	3	14,14,15	0.49	0	17,19,21	0.74	0
8	NAG	S	1	1,8	14,14,15	0.61	0	17,19,21	1.12	2 (11%)
8	NAG	S	2	8	14,14,15	0.56	0	17,19,21	0.95	1 (5%)
8	BMA	S	3	8	11,11,12	0.56	0	15,15,17	1.36	3 (20%)
8	MAN	S	4	8	11,11,12	0.53	0	15,15,17	1.02	1 (6%)
8	MAN	S	5	8	11,11,12	0.59	0	15,15,17	2.35	3 (20%)
8	FUC	S	6	8	10,10,11	0.76	0	14,14,16	2.78	5 (35%)
6	NAG	T	1	1,6	14,14,15	0.64	0	17,19,21	1.42	3 (17%)
6	NAG	T	2	6	14,14,15	0.61	0	17,19,21	0.99	1 (5%)
6	BMA	T	3	6	11,11,12	0.38	0	15,15,17	1.01	2 (13%)
9	NAG	U	1	9,2	14,14,15	0.55	0	17,19,21	1.30	3 (17%)
9	NAG	U	2	9	14,14,15	0.53	0	17,19,21	1.05	1 (5%)
9	FUC	U	3	9	10,10,11	0.78	0	14,14,16	1.41	2 (14%)
3	NAG	V	1	1,3	14,14,15	0.73	0	17,19,21	2.47	6 (35%)
3	NAG	V	2	3	14,14,15	0.46	0	17,19,21	1.96	2 (11%)
3	NAG	W	1	1,3	14,14,15	0.66	0	17,19,21	1.67	3 (17%)
3	NAG	W	2	3	14,14,15	0.53	0	17,19,21	0.91	1 (5%)
4	NAG	X	1	1,4	14,14,15	0.63	0	17,19,21	0.94	1 (5%)
4	NAG	X	2	4	14,14,15	0.65	0	17,19,21	0.91	1 (5%)
4	BMA	X	3	4	11,11,12	0.37	0	15,15,17	1.22	1 (6%)
4	MAN	X	4	4	11,11,12	0.60	0	15,15,17	0.95	1 (6%)
4	MAN	X	5	4	11,11,12	0.62	0	15,15,17	1.16	1 (6%)
3	NAG	Y	1	1,3	14,14,15	0.62	0	17,19,21	1.80	4 (23%)
3	NAG	Y	2	3	14,14,15	0.62	0	17,19,21	2.15	4 (23%)
9	NAG	Z	1	1,9	14,14,15	0.51	0	17,19,21	1.14	1 (5%)
9	NAG	Z	2	9	14,14,15	0.64	0	17,19,21	0.93	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	FUC	Z	3	9	10,10,11	0.81	0	14,14,16	1.73	3 (21%)
3	NAG	a	1	1,3	14,14,15	0.64	0	17,19,21	1.35	3 (17%)
3	NAG	a	2	3	14,14,15	0.54	0	17,19,21	0.93	1 (5%)
9	NAG	b	1	9,2	14,14,15	0.56	0	17,19,21	1.61	3 (17%)
9	NAG	b	2	9	14,14,15	0.42	0	17,19,21	1.04	1 (5%)
9	FUC	b	3	9	10,10,11	0.73	0	14,14,16	0.80	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
3	NAG	G	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	G	2	3	-	4/6/23/26	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
4	NAG	I	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	0/6/23/26	0/1/1/1
4	BMA	I	3	4	-	2/2/19/22	0/1/1/1
4	MAN	I	4	4	-	2/2/19/22	0/1/1/1
4	MAN	I	5	4	-	2/2/19/22	0/1/1/1
3	NAG	J	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	J	2	3	-	1/6/23/26	0/1/1/1
5	NAG	K	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	K	2	5	-	1/6/23/26	0/1/1/1
5	BMA	K	3	5	-	2/2/19/22	0/1/1/1
5	FUC	K	4	5	1/1/4/5	-	0/1/1/1
3	NAG	L	1	1,3	1/1/5/7	3/6/23/26	0/1/1/1
3	NAG	L	2	3	-	2/6/23/26	0/1/1/1
6	NAG	M	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	M	2	6	-	2/6/23/26	0/1/1/1
6	BMA	M	3	6	-	2/2/19/22	0/1/1/1
7	NAG	N	1	2,7	-	2/6/23/26	0/1/1/1
7	FUC	N	2	7	1/1/4/5	-	0/1/1/1
6	NAG	O	1	1,6	-	1/6/23/26	0/1/1/1
6	NAG	O	2	6	-	2/6/23/26	0/1/1/1
6	BMA	O	3	6	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	P	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	P	2	4	-	0/6/23/26	0/1/1/1
4	BMA	P	3	4	-	2/2/19/22	0/1/1/1
4	MAN	P	4	4	-	2/2/19/22	0/1/1/1
4	MAN	P	5	4	-	2/2/19/22	0/1/1/1
4	NAG	Q	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	Q	2	4	-	0/6/23/26	0/1/1/1
4	BMA	Q	3	4	-	0/2/19/22	0/1/1/1
4	MAN	Q	4	4	-	0/2/19/22	1/1/1/1
4	MAN	Q	5	4	-	2/2/19/22	0/1/1/1
3	NAG	R	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	R	2	3	-	0/6/23/26	0/1/1/1
8	NAG	S	1	1,8	-	2/6/23/26	0/1/1/1
8	NAG	S	2	8	-	0/6/23/26	0/1/1/1
8	BMA	S	3	8	-	2/2/19/22	0/1/1/1
8	MAN	S	4	8	-	0/2/19/22	0/1/1/1
8	MAN	S	5	8	-	1/2/19/22	0/1/1/1
8	FUC	S	6	8	1/1/4/5	-	0/1/1/1
6	NAG	T	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	T	2	6	-	0/6/23/26	0/1/1/1
6	BMA	T	3	6	-	2/2/19/22	0/1/1/1
9	NAG	U	1	9,2	-	2/6/23/26	0/1/1/1
9	NAG	U	2	9	-	0/6/23/26	0/1/1/1
9	FUC	U	3	9	1/1/4/5	-	0/1/1/1
3	NAG	V	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	V	2	3	-	2/6/23/26	0/1/1/1
3	NAG	W	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	W	2	3	-	0/6/23/26	0/1/1/1
4	NAG	X	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	X	2	4	-	0/6/23/26	0/1/1/1
4	BMA	X	3	4	-	0/2/19/22	0/1/1/1
4	MAN	X	4	4	-	2/2/19/22	0/1/1/1
4	MAN	X	5	4	-	2/2/19/22	0/1/1/1
3	NAG	Y	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	Y	2	3	-	6/6/23/26	0/1/1/1
9	NAG	Z	1	1,9	-	2/6/23/26	0/1/1/1
9	NAG	Z	2	9	-	0/6/23/26	0/1/1/1
9	FUC	Z	3	9	1/1/4/5	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	a	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	a	2	3	-	2/6/23/26	0/1/1/1
9	NAG	b	1	9,2	-	2/6/23/26	0/1/1/1
9	NAG	b	2	9	-	2/6/23/26	0/1/1/1
9	FUC	b	3	9	1/1/4/5	-	0/1/1/1

There are no bond length outliers.

The worst 5 of 123 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	S	6	FUC	C1-C2-C3	-8.04	99.78	109.67
3	V	2	NAG	C1-O5-C5	7.02	121.71	112.19
3	Y	2	NAG	C2-N2-C7	6.60	132.30	122.90
3	V	1	NAG	C1-O5-C5	6.38	120.84	112.19
4	I	3	BMA	C1-C2-C3	6.32	117.43	109.67

5 of 7 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
9	Z	3	FUC	C1
9	U	3	FUC	C1
7	N	2	FUC	C1
3	L	1	NAG	C1
9	b	3	FUC	C1

5 of 90 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	T	3	BMA	O5-C5-C6-O6
4	X	5	MAN	O5-C5-C6-O6
8	S	1	NAG	C4-C5-C6-O6
9	b	1	NAG	O5-C5-C6-O6
6	O	3	BMA	O5-C5-C6-O6

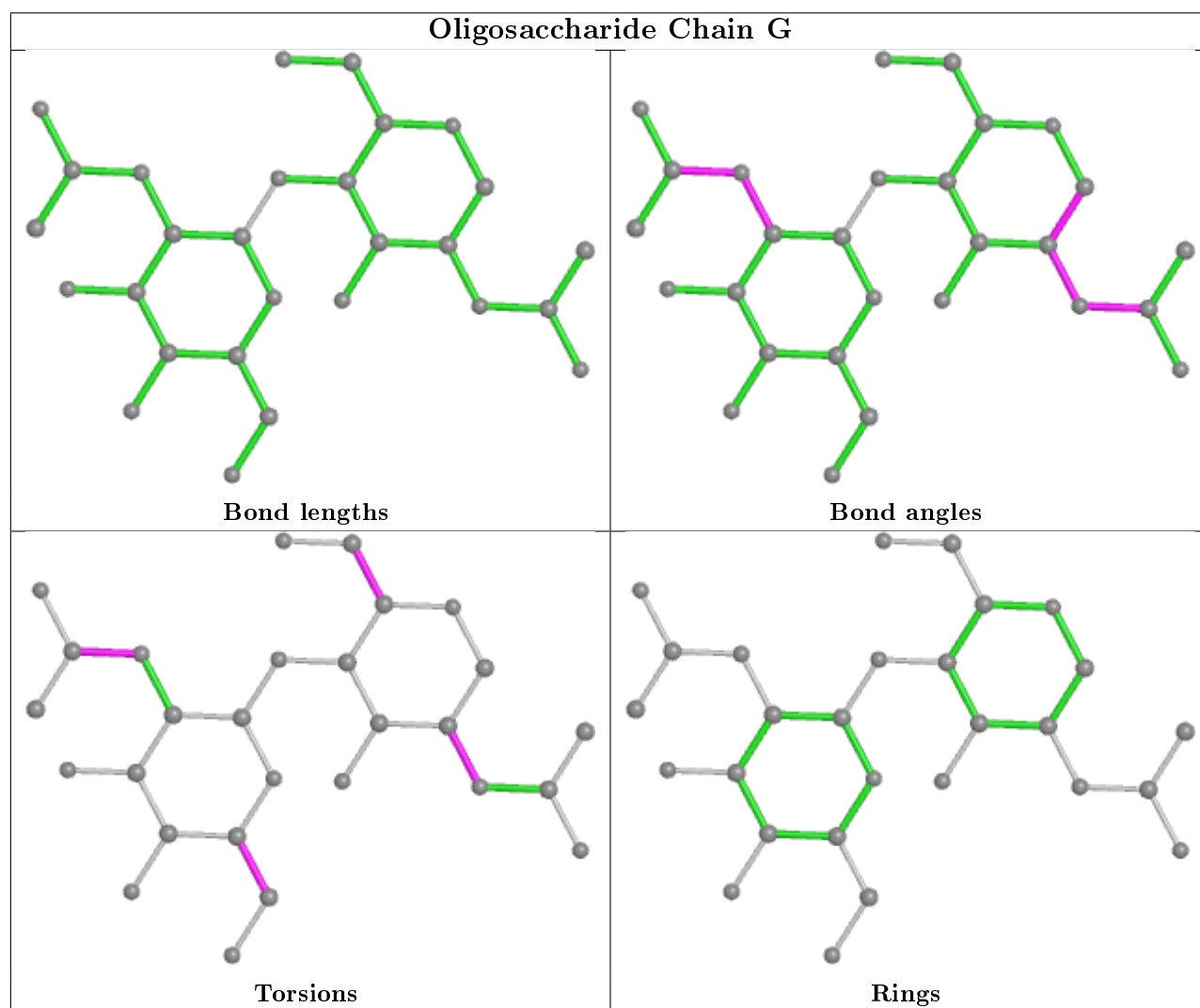
All (1) ring outliers are listed below:

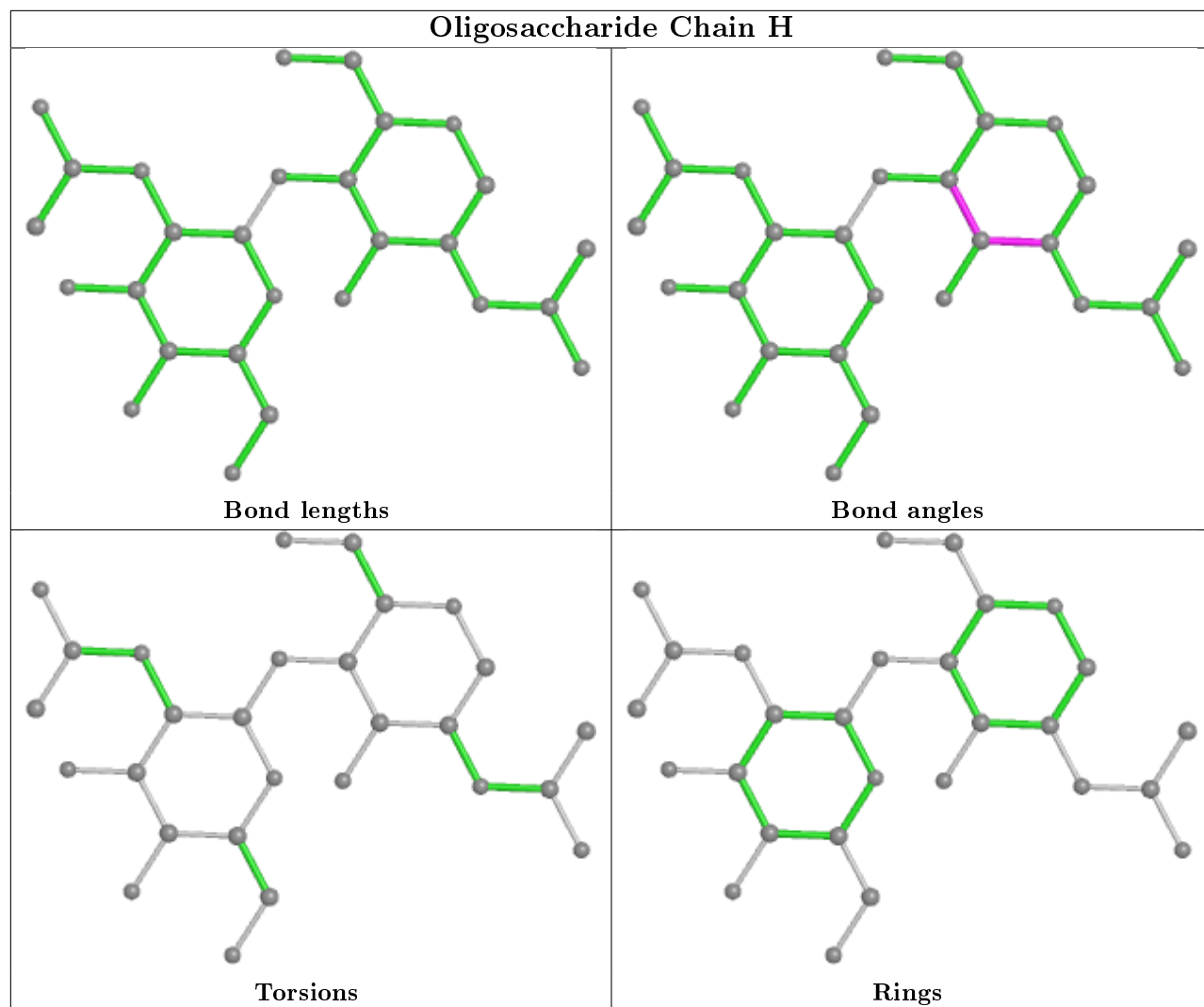
Mol	Chain	Res	Type	Atoms
4	Q	4	MAN	C1-C2-C3-C4-C5-O5

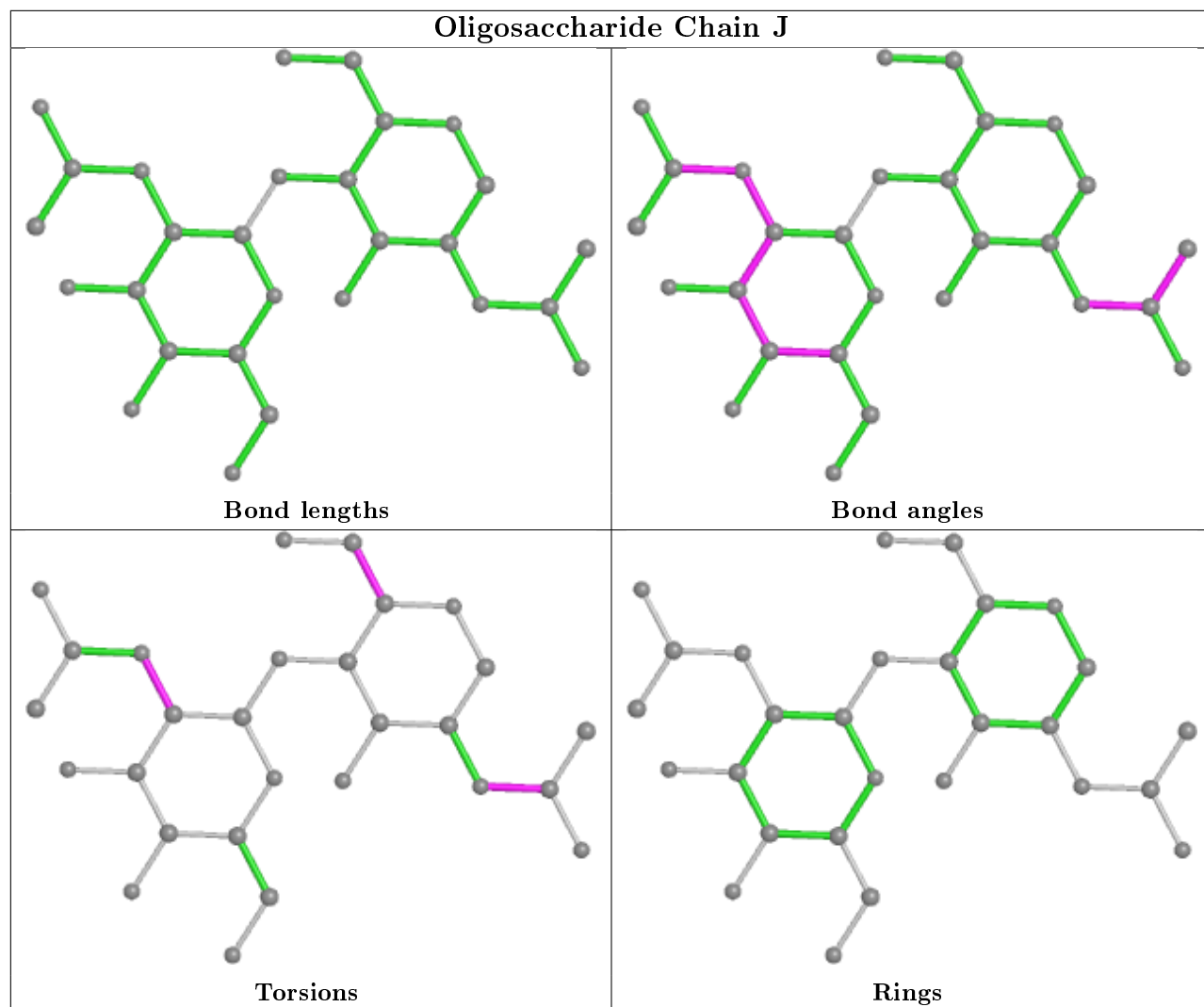
11 monomers are involved in 12 short contacts:

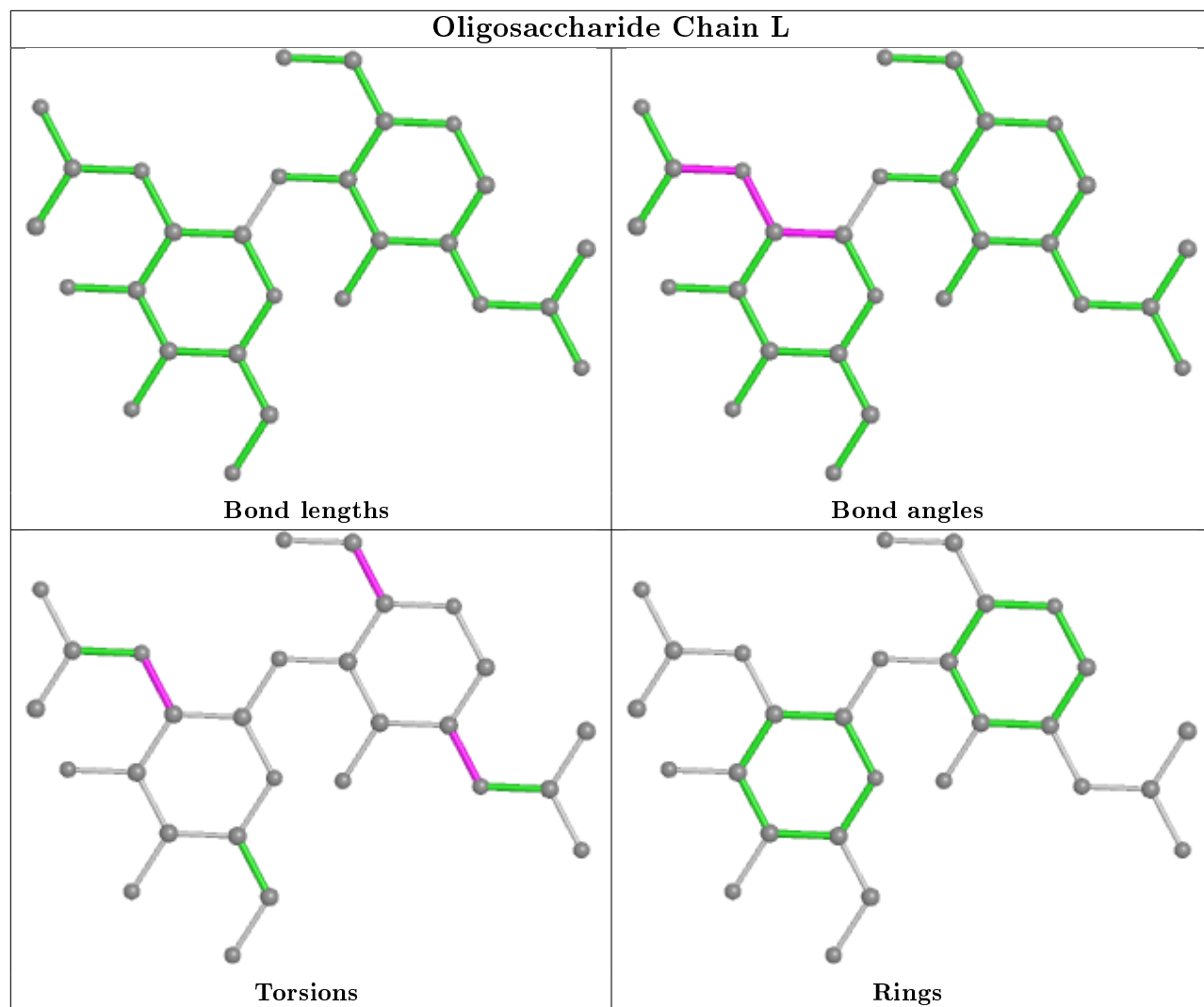
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	P	2	NAG	3	0
4	P	3	BMA	4	0
4	P	4	MAN	1	0
4	Q	1	NAG	1	0
3	V	1	NAG	1	0
6	T	1	NAG	1	0
4	I	5	MAN	2	0
3	V	2	NAG	1	0
3	H	1	NAG	1	0
4	I	3	BMA	2	0
5	K	4	FUC	2	0

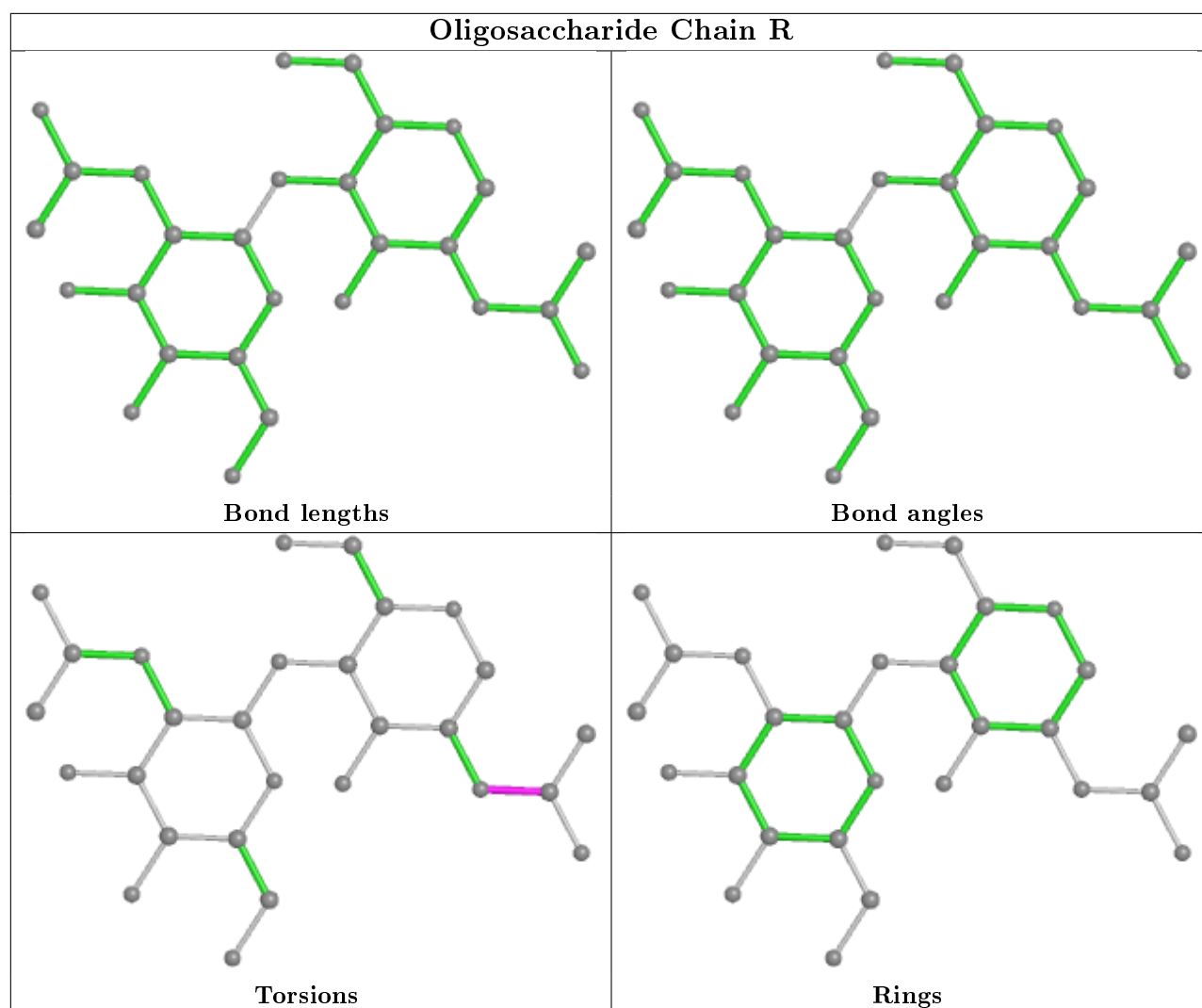
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

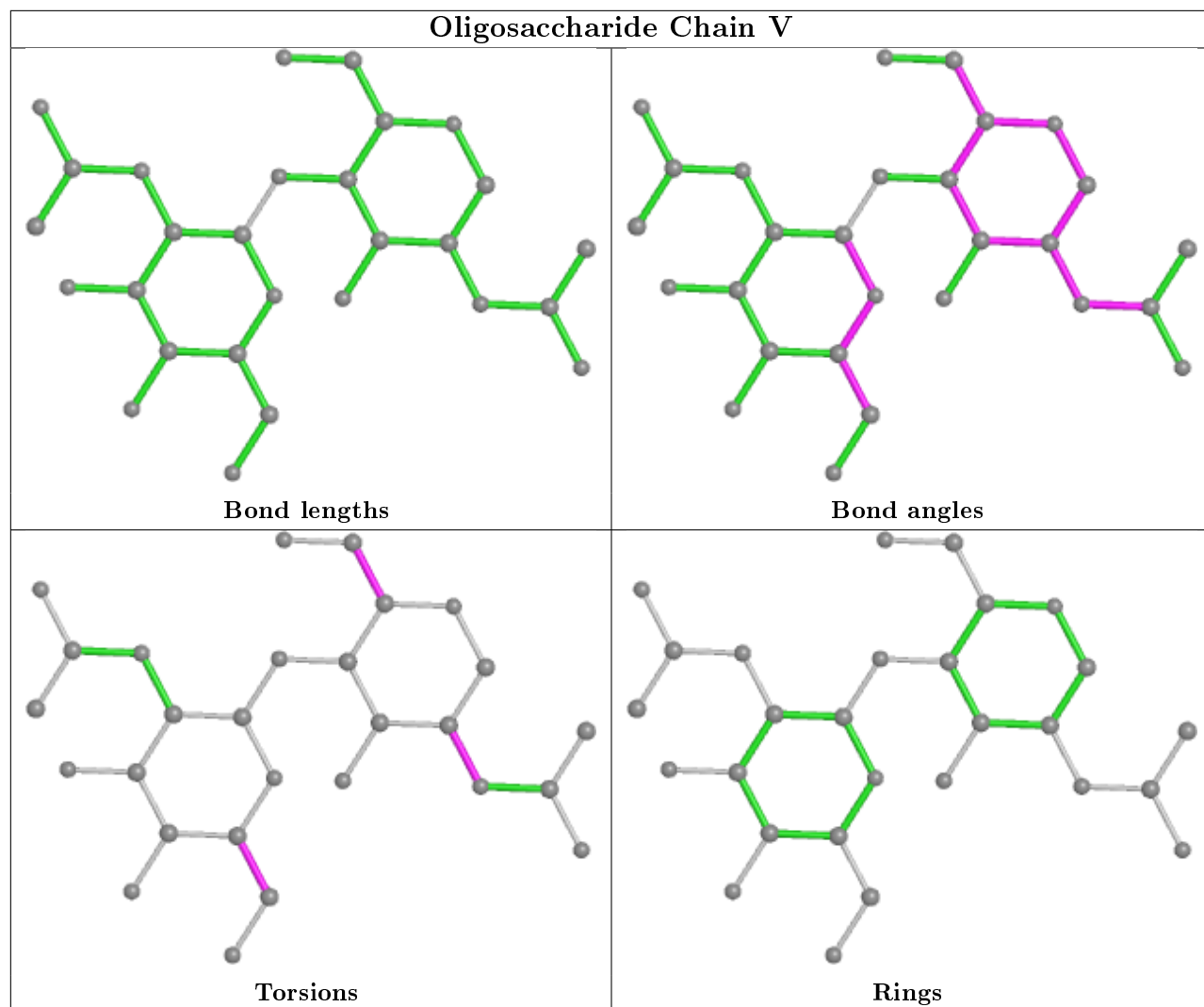


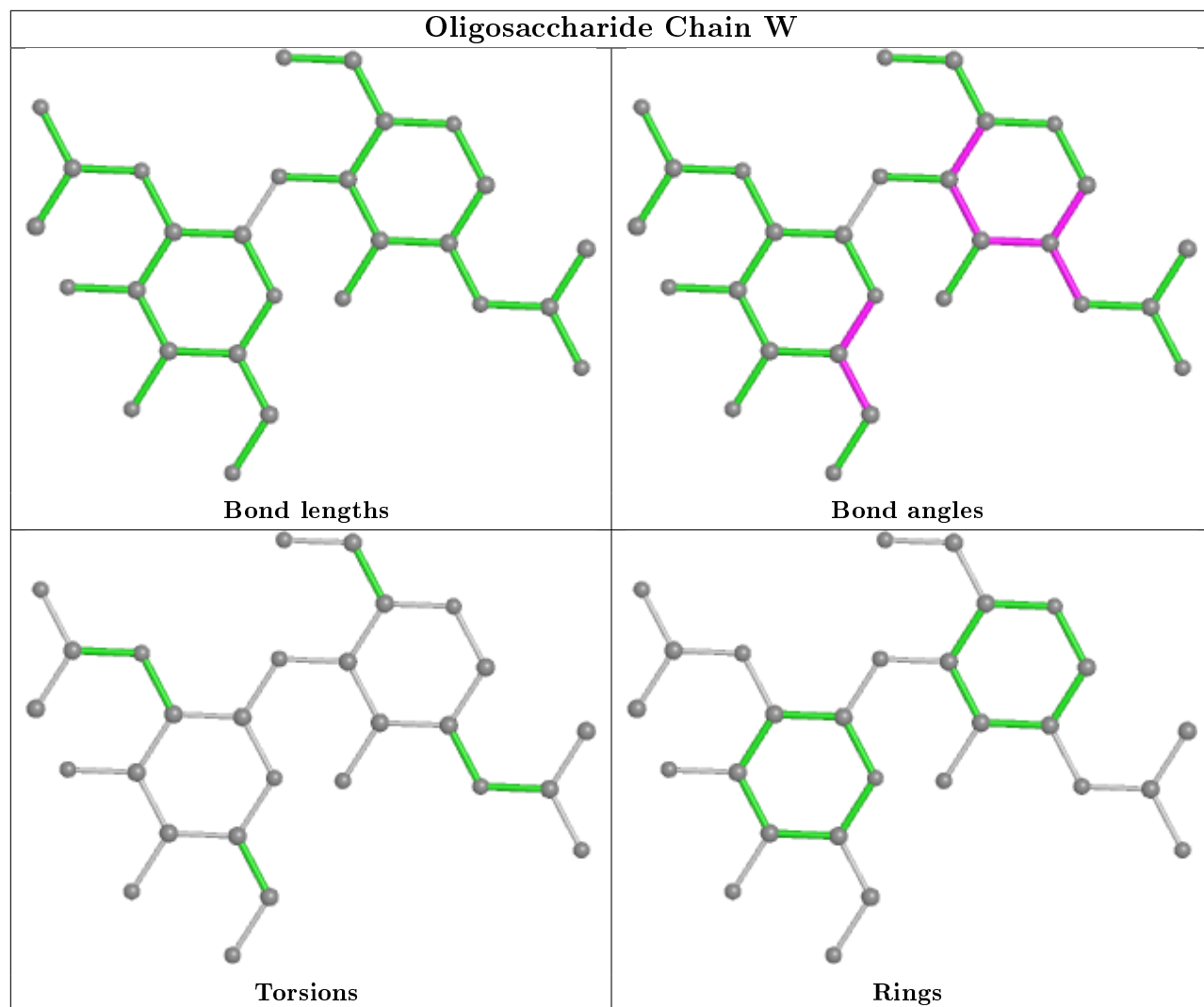


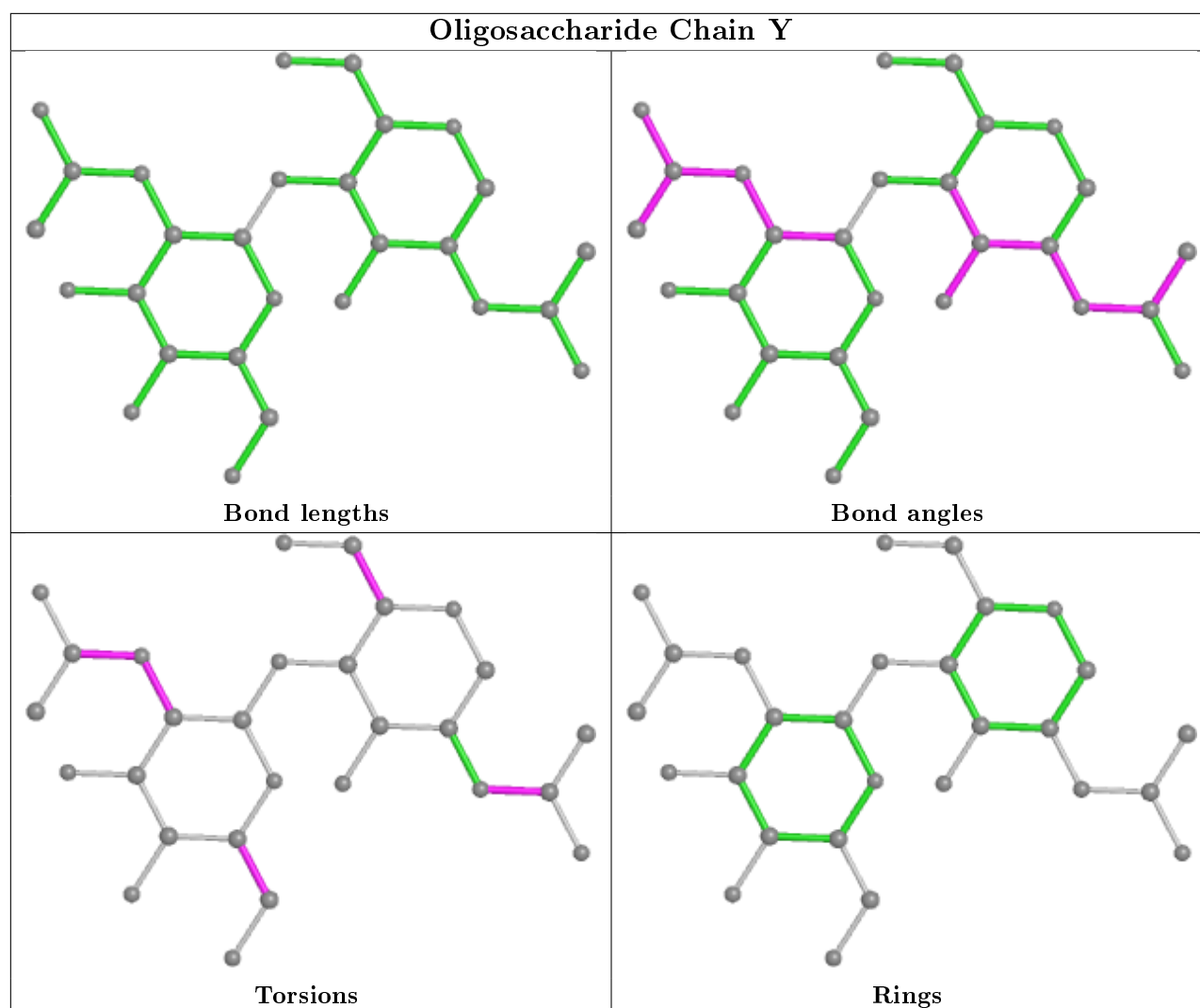


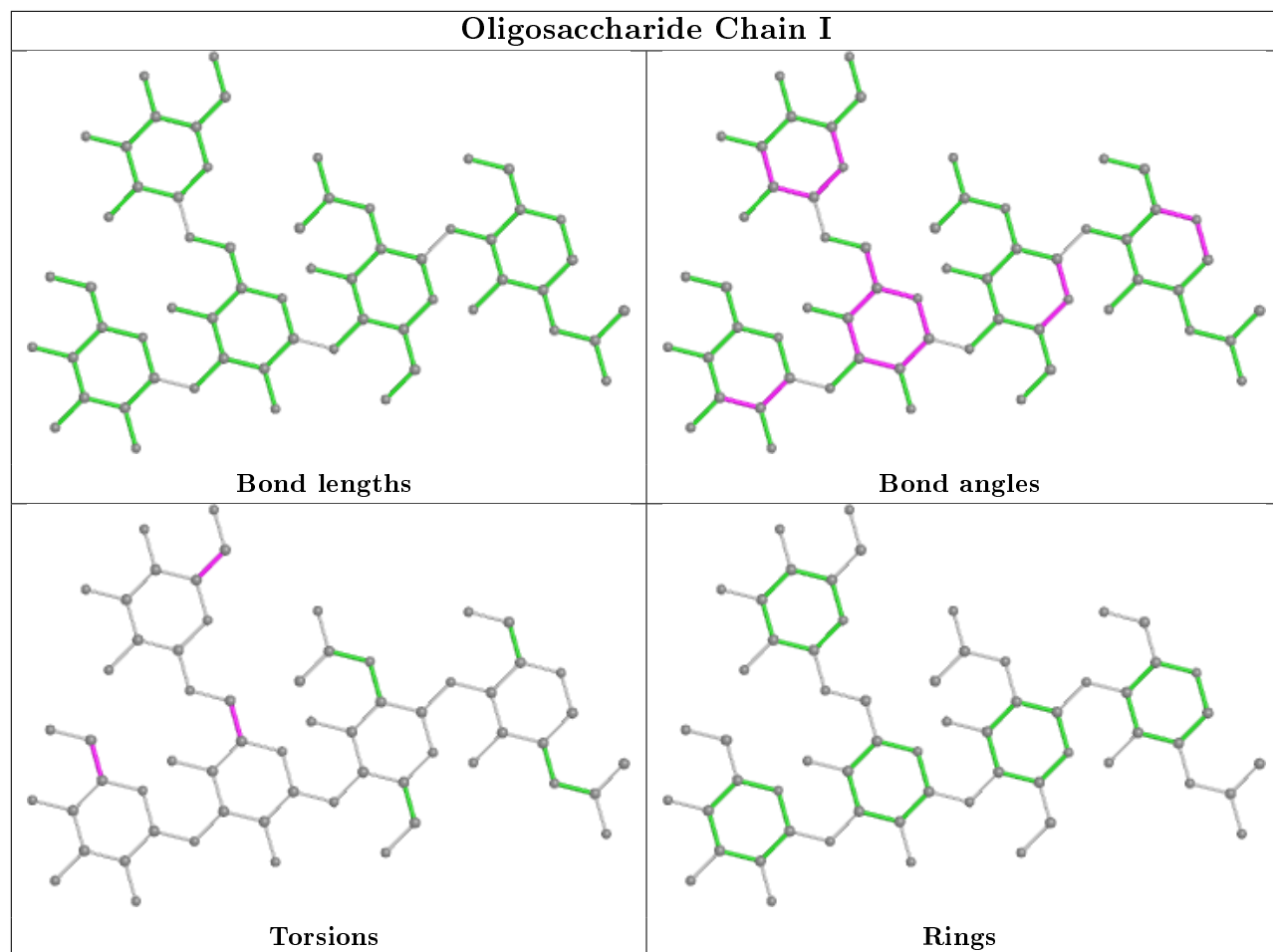


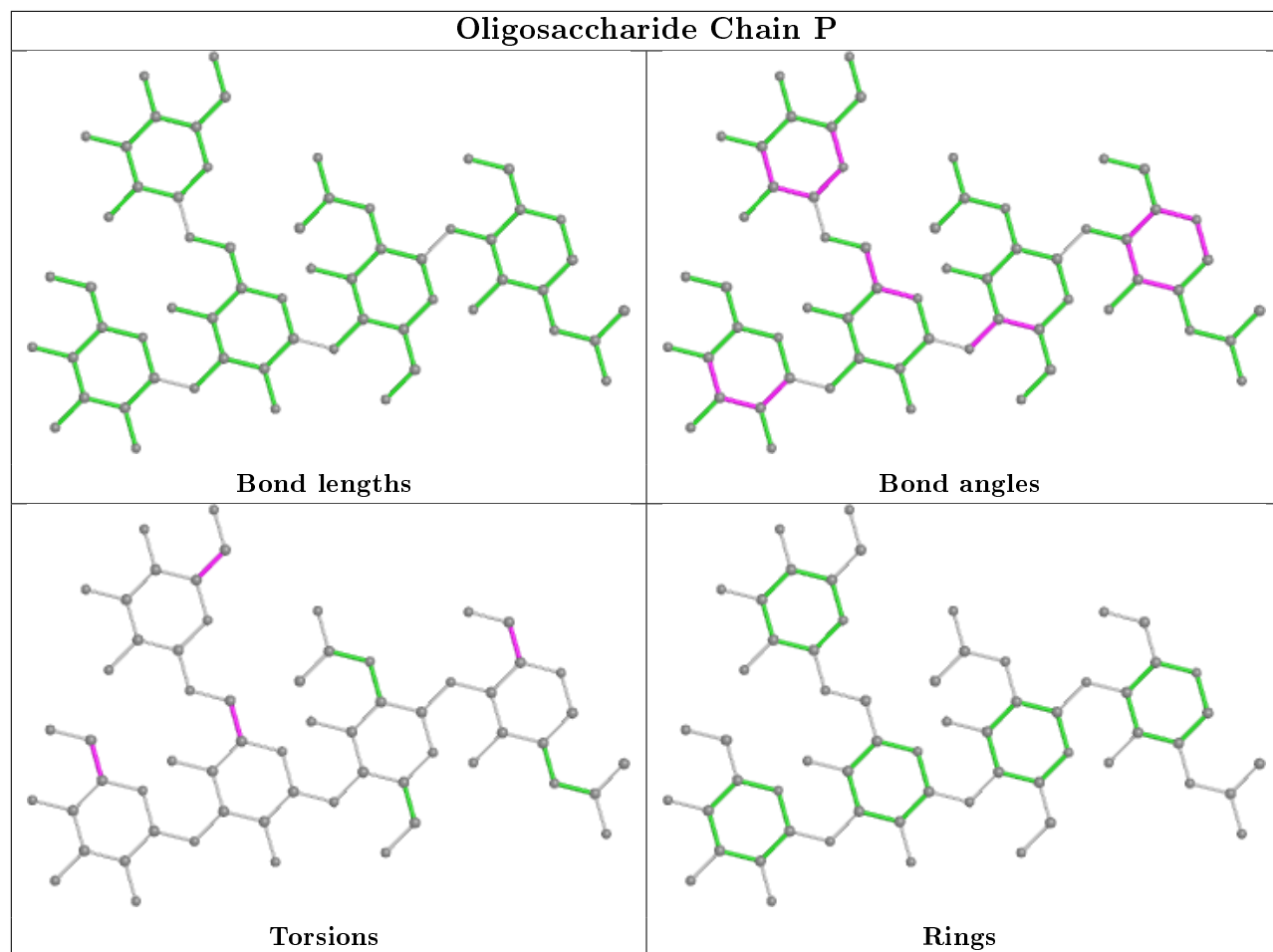


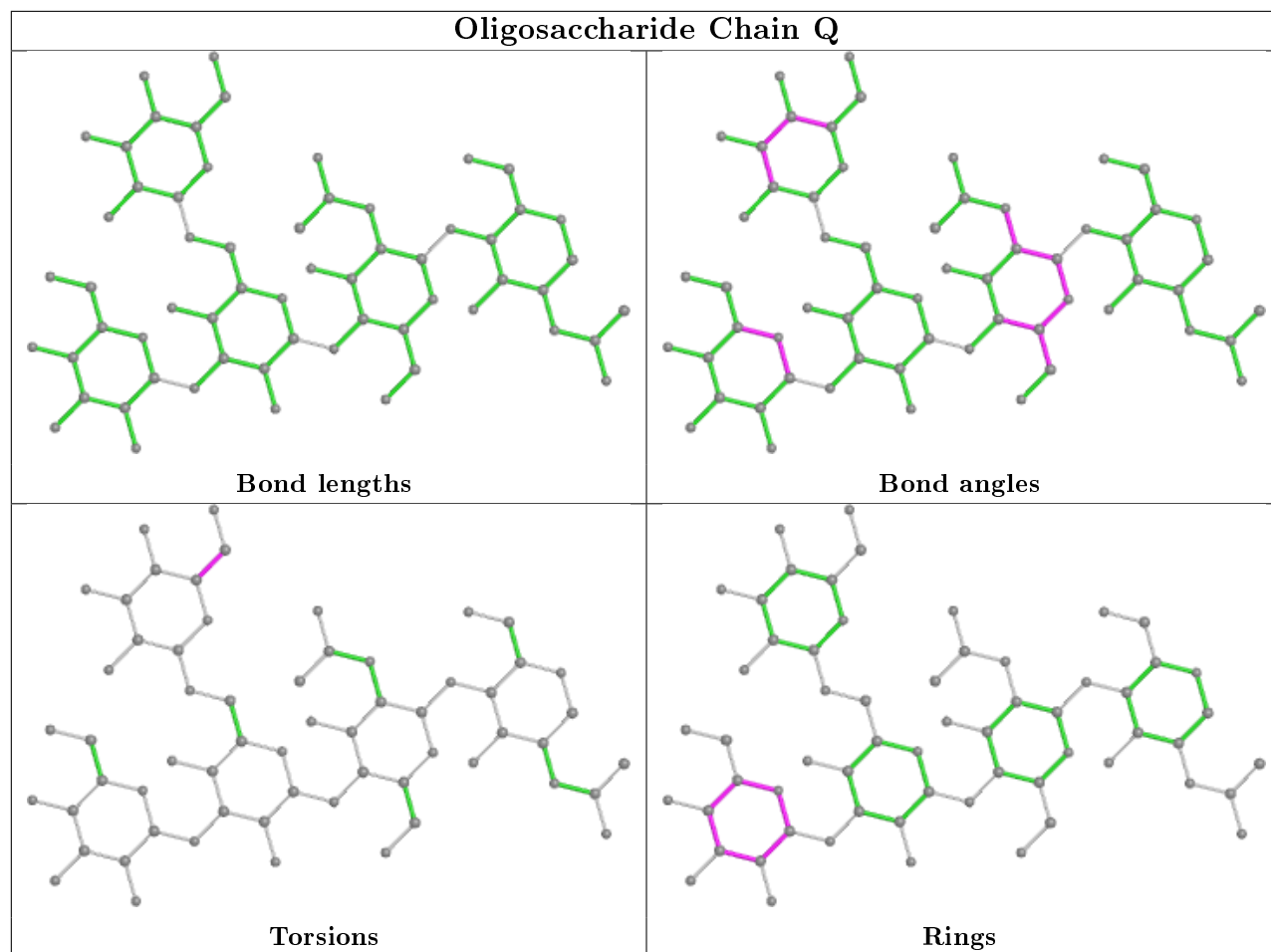


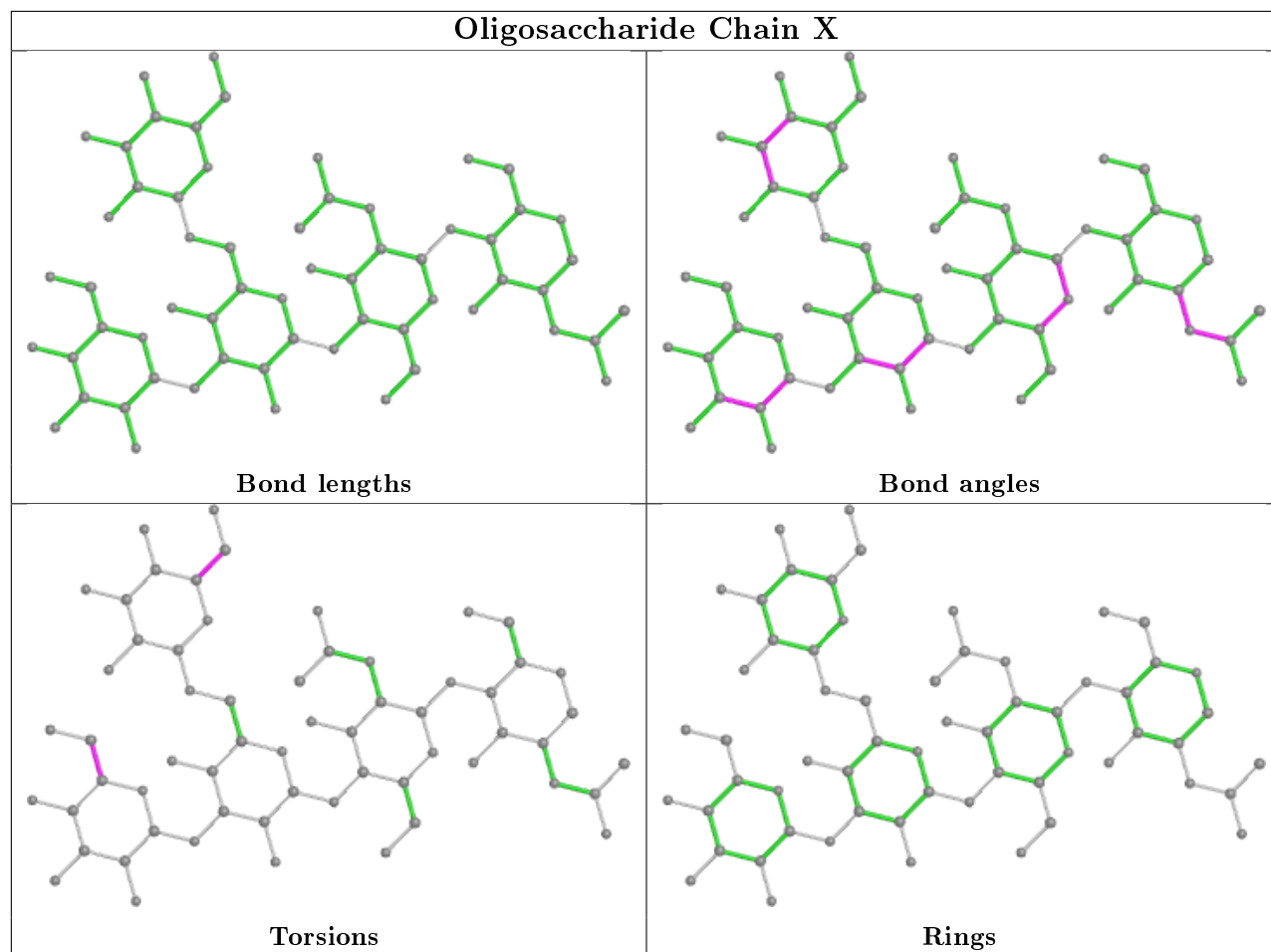


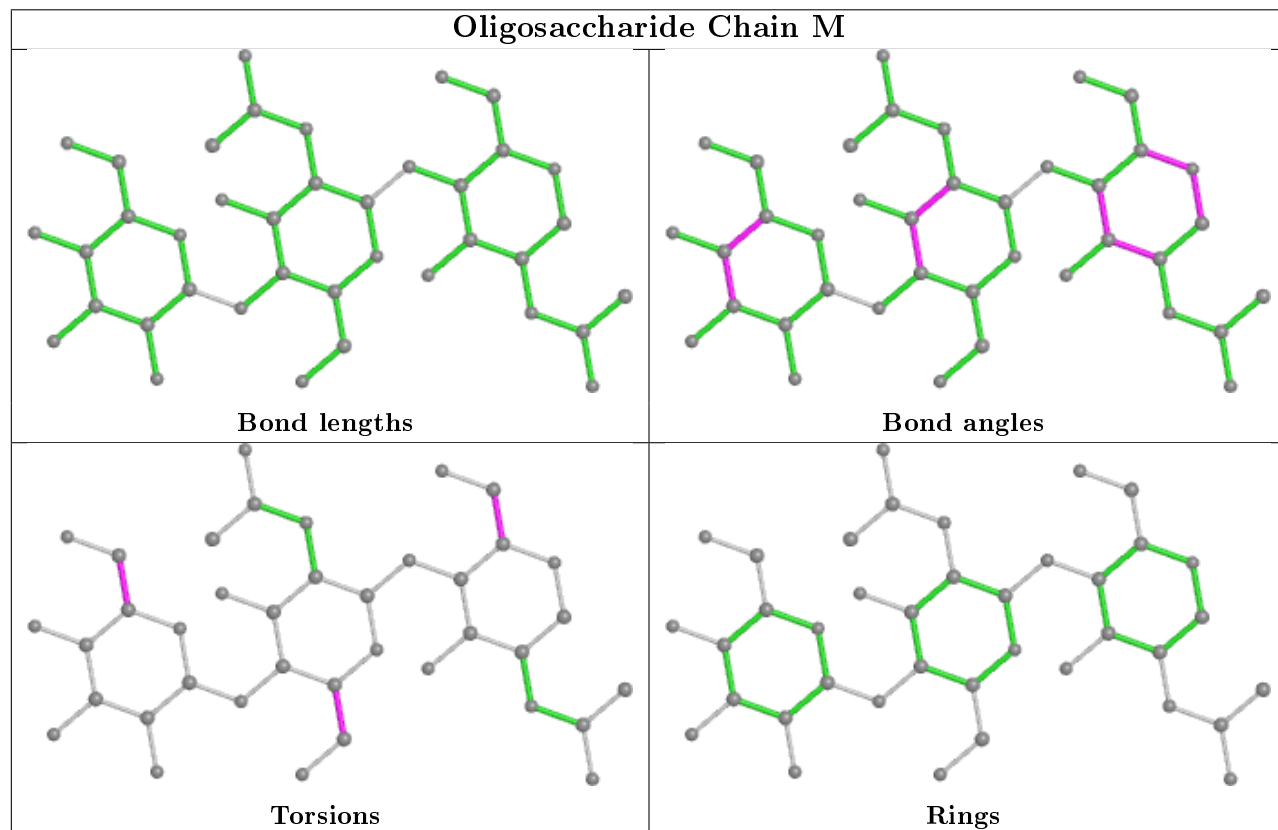
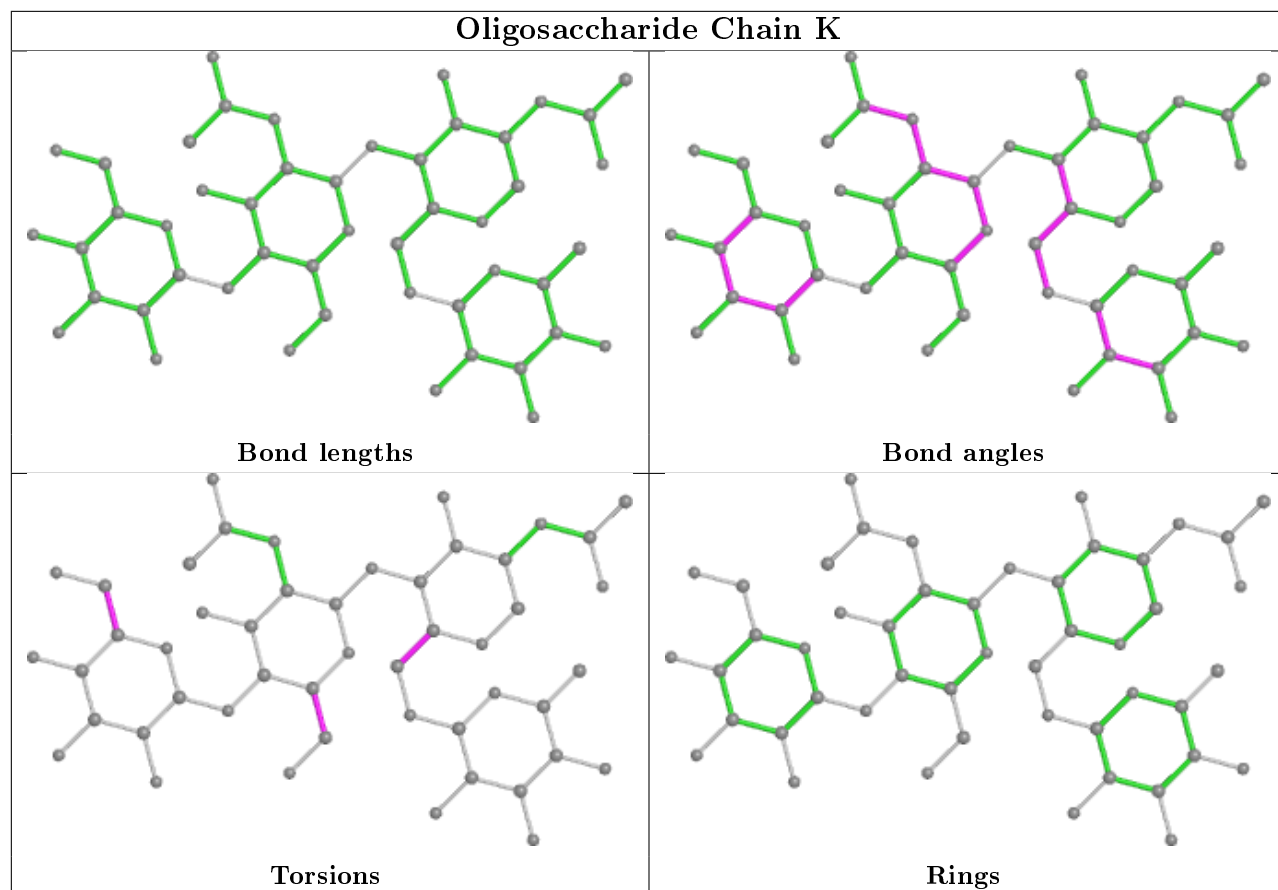


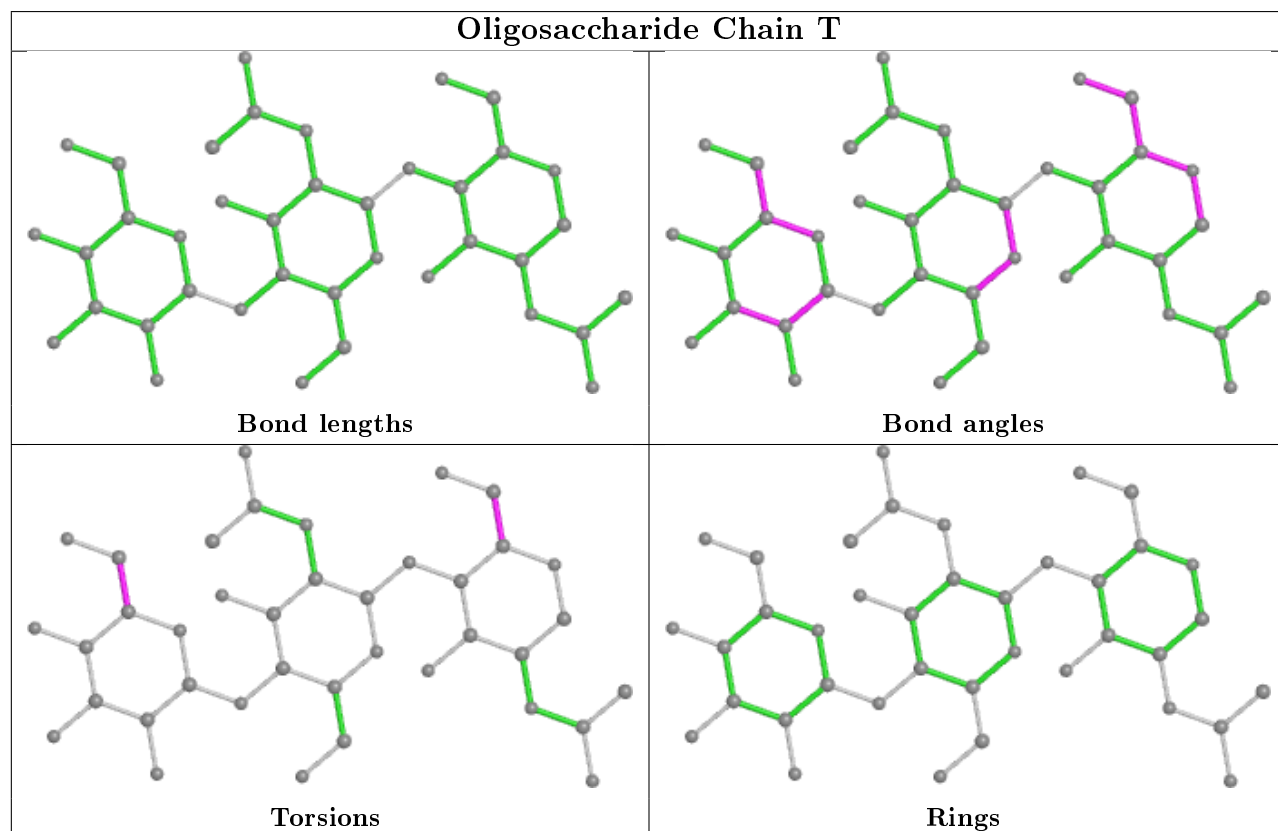
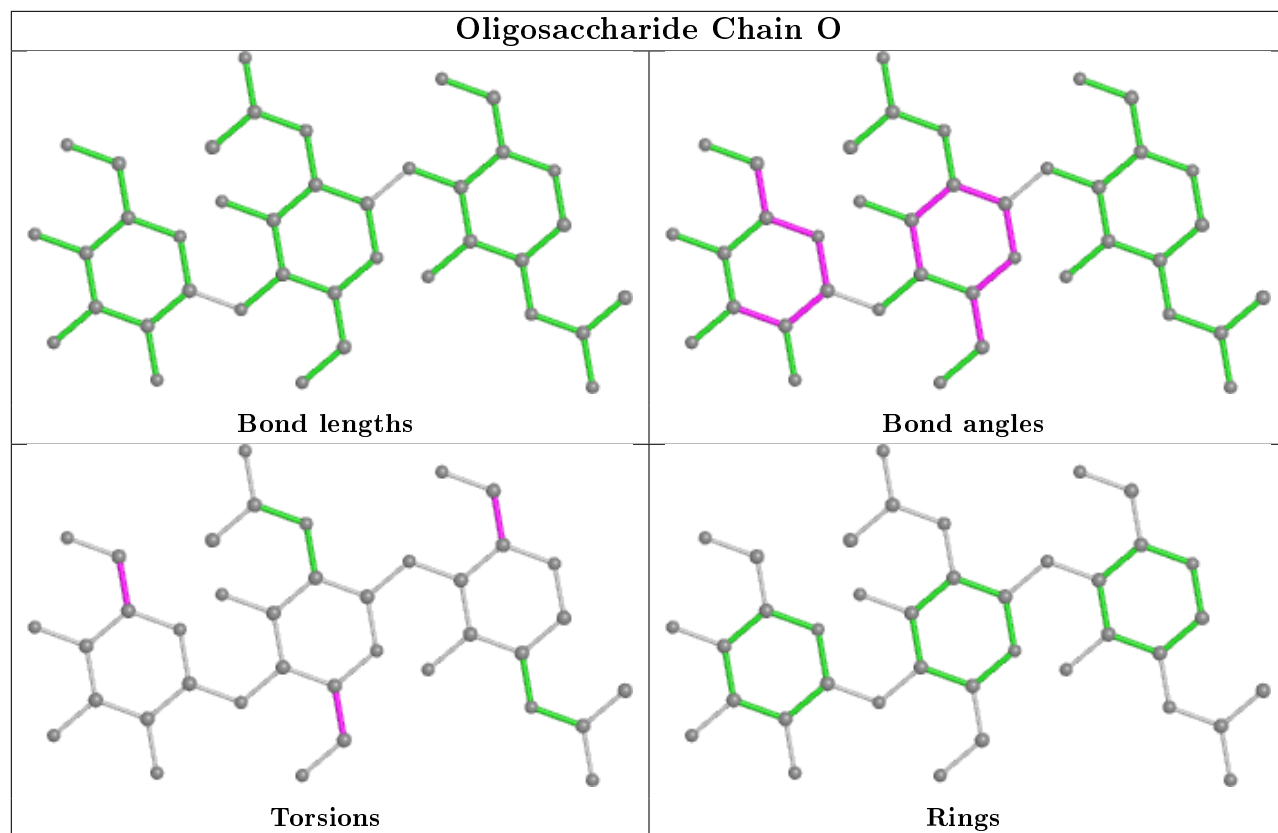


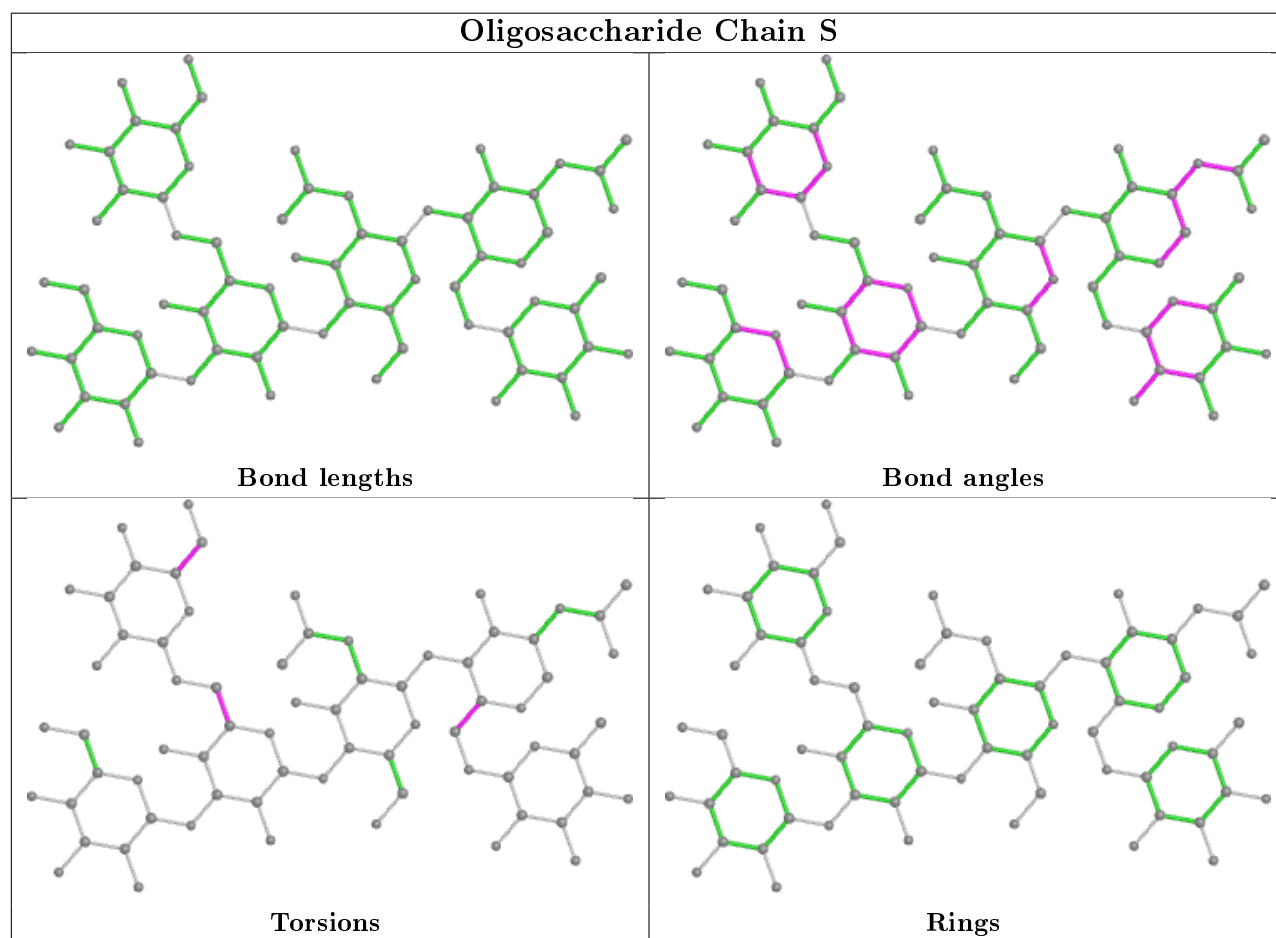
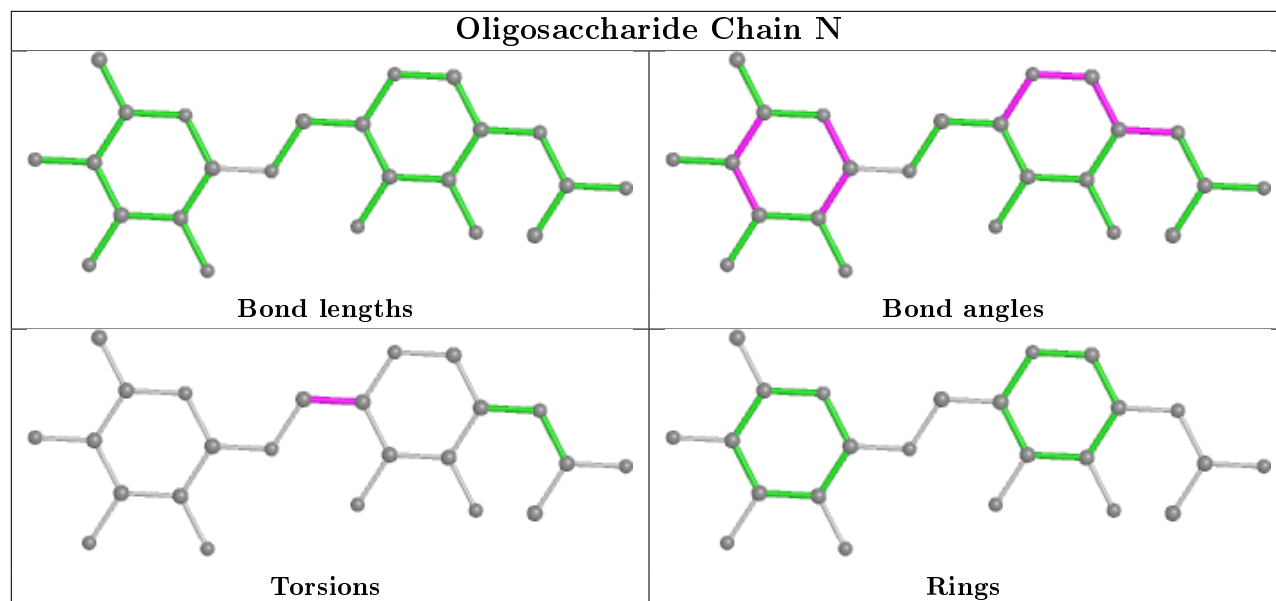


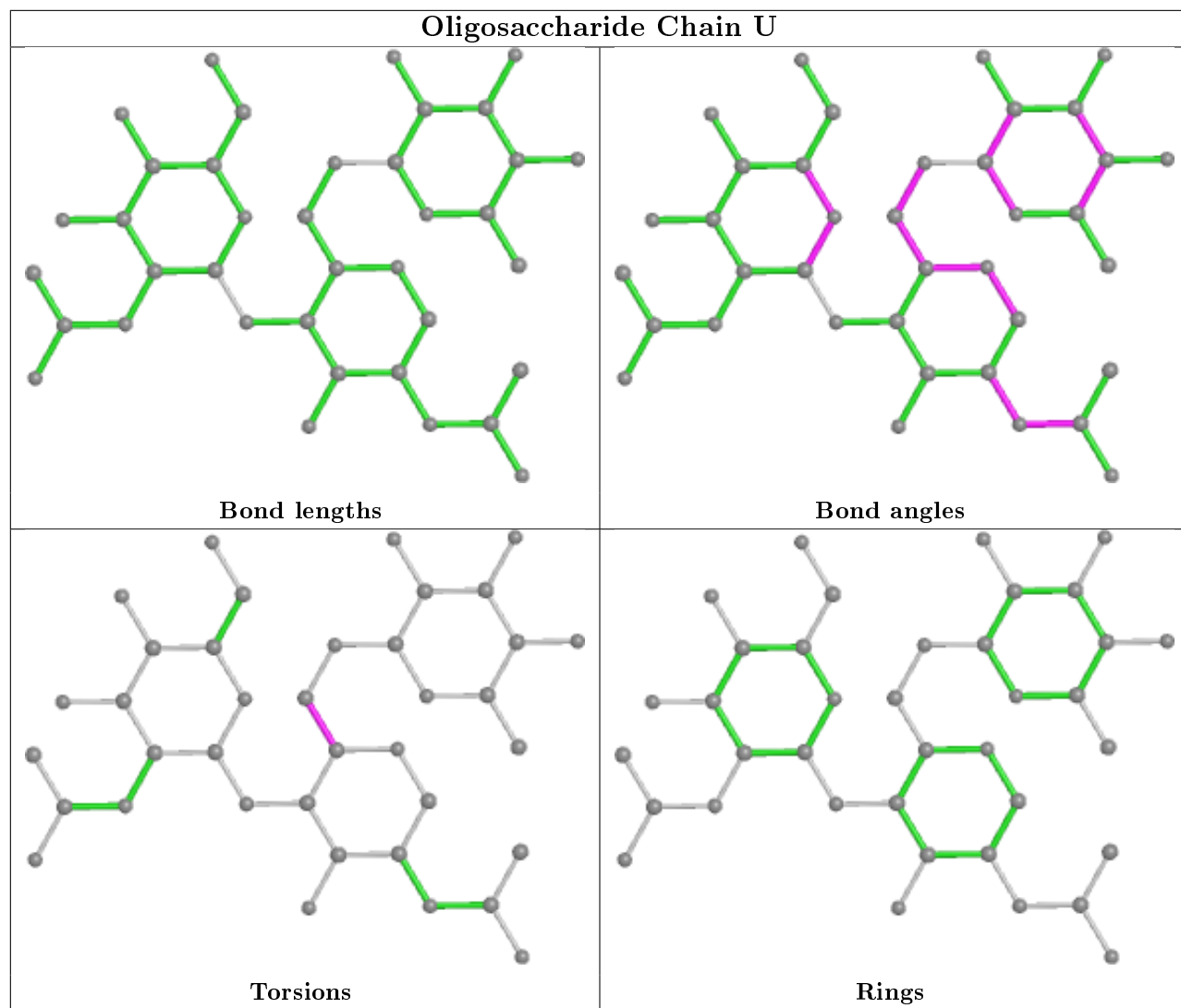


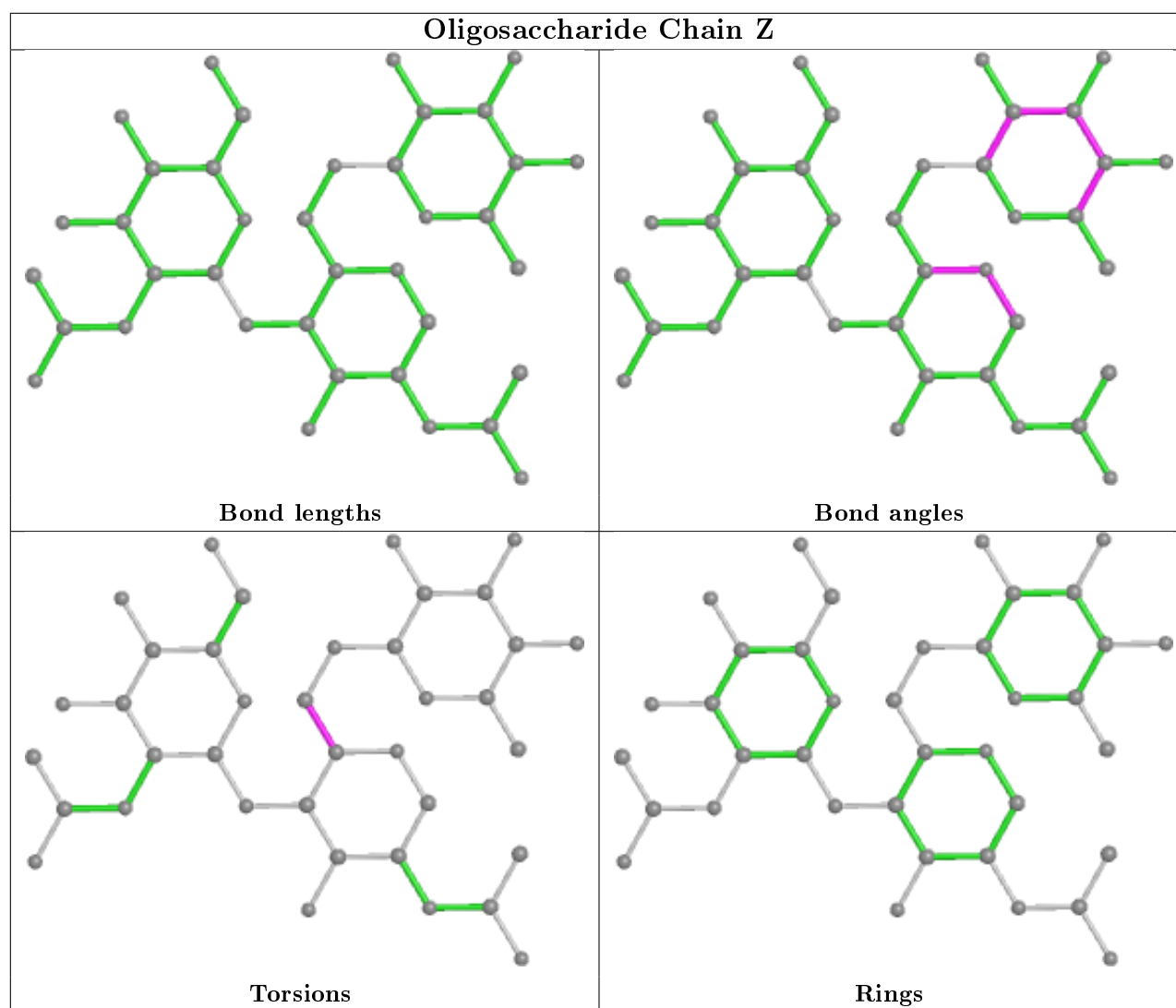












5.6 Ligand geometry [i](#)

15 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	$\# Z > 2$
10	BMA	A	604	-	11,11,12	0.38	0	15,15,17	0.72	0
12	NAG	C	601	1	14,14,15	0.54	0	17,19,21	1.01	1 (5%)
11	EDO	E	1326	-	3,3,3	0.45	0	2,2,2	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	EDO	A	1328	-	3,3,3	0.47	0	2,2,2	0.37	0
11	EDO	F	1174	-	3,3,3	0.51	0	2,2,2	0.15	0
11	EDO	B	1174	-	3,3,3	0.46	0	2,2,2	0.36	0
11	EDO	C	1328	-	3,3,3	0.47	0	2,2,2	0.45	0
11	EDO	B	1173	-	3,3,3	0.38	0	2,2,2	0.50	0
11	EDO	D	1173	-	3,3,3	0.46	0	2,2,2	0.37	0
12	NAG	E	621	1	14,14,15	0.65	0	17,19,21	1.11	1 (5%)
11	EDO	F	1173	-	3,3,3	0.46	0	2,2,2	0.41	0
11	EDO	A	1327	-	3,3,3	0.35	0	2,2,2	0.62	0
11	EDO	C	1327	-	3,3,3	0.48	0	2,2,2	0.37	0
11	EDO	E	1327	-	3,3,3	0.47	0	2,2,2	0.38	0
11	EDO	F	1175	-	3,3,3	0.46	0	2,2,2	0.38	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
10	BMA	A	604	-	-	2/2/19/22	0/1/1/1
12	NAG	C	601	1	-	4/6/23/26	0/1/1/1
11	EDO	E	1326	-	-	1/1/1/1	-
11	EDO	A	1328	-	-	1/1/1/1	-
11	EDO	F	1174	-	-	1/1/1/1	-
11	EDO	B	1174	-	-	0/1/1/1	-
11	EDO	C	1328	-	-	0/1/1/1	-
11	EDO	B	1173	-	-	0/1/1/1	-
11	EDO	D	1173	-	-	1/1/1/1	-
12	NAG	E	621	1	-	2/6/23/26	0/1/1/1
11	EDO	F	1173	-	-	1/1/1/1	-
11	EDO	A	1327	-	-	1/1/1/1	-
11	EDO	C	1327	-	-	0/1/1/1	-
11	EDO	E	1327	-	-	1/1/1/1	-
11	EDO	F	1175	-	-	1/1/1/1	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	E	621	NAG	C4-C3-C2	3.60	116.29	111.02
12	C	601	NAG	C1-O5-C5	2.03	114.94	112.19

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	C	601	NAG	O5-C5-C6-O6
12	E	621	NAG	C4-C5-C6-O6
12	C	601	NAG	C8-C7-N2-C2
12	C	601	NAG	O7-C7-N2-C2
12	E	621	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	A	1328	EDO	1	0
11	E	1327	EDO	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	325/329 (98%)	-0.04	6 (1%) 68 71	19, 29, 48, 64	0
1	C	320/329 (97%)	-0.03	2 (0%) 89 90	17, 29, 48, 71	0
1	E	325/329 (98%)	-0.07	3 (0%) 84 85	18, 30, 47, 66	0
2	B	172/172 (100%)	0.03	2 (1%) 79 81	15, 26, 42, 59	0
2	D	172/172 (100%)	0.10	4 (2%) 60 63	15, 27, 44, 63	0
2	F	172/172 (100%)	0.17	2 (1%) 79 81	16, 24, 41, 61	0
All	All	1486/1503 (98%)	0.00	19 (1%) 77 79	15, 28, 47, 71	0

The worst 5 of 19 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	81	TYR	4.7
1	A	276	ILE	4.2
1	A	142	GLY	3.7
1	A	278	VAL	3.3
2	D	58	ARG	3.2

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 [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	MAN	P	5	11/12	0.50	0.56	94,100,104,107	0
3	NAG	L	1	14/15	0.55	0.34	66,72,76,82	0
6	BMA	O	3	11/12	0.56	0.30	90,95,101,101	0
3	NAG	W	2	14/15	0.56	0.40	85,89,96,97	0
3	NAG	J	2	14/15	0.58	0.36	71,79,82,84	0
6	BMA	M	3	11/12	0.59	0.28	84,87,91,93	0
3	NAG	V	2	14/15	0.59	0.39	95,99,101,101	0
3	NAG	Y	2	14/15	0.59	0.47	81,84,87,90	0
3	NAG	G	2	14/15	0.63	0.49	94,100,105,105	0
3	NAG	H	2	14/15	0.65	0.50	84,88,94,95	0
4	MAN	I	4	11/12	0.66	0.22	79,86,93,95	0
4	MAN	P	4	11/12	0.68	0.47	84,87,90,90	0
3	NAG	R	2	14/15	0.69	0.31	72,77,80,80	0
9	NAG	b	2	14/15	0.69	0.34	82,87,91,93	0
9	NAG	U	2	14/15	0.69	0.35	74,80,85,85	0
9	NAG	Z	2	14/15	0.71	0.19	70,72,77,78	0
4	MAN	Q	5	11/12	0.71	0.27	79,80,84,85	0
3	NAG	W	1	14/15	0.73	0.30	61,69,74,80	0
6	NAG	O	2	14/15	0.73	0.35	82,85,89,89	0
3	NAG	L	2	14/15	0.74	0.39	83,89,93,94	0
3	NAG	J	1	14/15	0.74	0.20	51,62,66,71	0
3	NAG	R	1	14/15	0.74	0.24	52,61,65,68	0
9	FUC	Z	3	10/11	0.74	0.28	71,76,79,79	0
4	MAN	Q	4	11/12	0.75	0.27	70,75,81,83	0
6	NAG	M	2	14/15	0.75	0.26	70,75,80,84	0
4	BMA	I	3	11/12	0.76	0.12	64,71,78,78	0
4	BMA	P	3	11/12	0.77	0.34	84,86,91,97	0
5	BMA	K	3	11/12	0.77	0.36	80,86,89,92	0
6	NAG	O	1	14/15	0.78	0.30	65,69,74,78	0
9	FUC	U	3	10/11	0.78	0.26	62,68,71,72	0
9	NAG	b	1	14/15	0.79	0.21	54,62,68,76	0
8	MAN	S	5	11/12	0.79	0.40	72,74,81,85	0
4	MAN	X	4	11/12	0.80	0.25	78,81,84,87	0
9	FUC	b	3	10/11	0.80	0.38	74,78,80,82	0
3	NAG	V	1	14/15	0.80	0.35	76,82,88,90	0
4	BMA	Q	3	11/12	0.81	0.15	62,69,74,74	0
4	MAN	X	5	11/12	0.81	0.20	77,80,84,85	0
5	NAG	K	2	14/15	0.82	0.26	58,63,71,75	0
4	NAG	P	2	14/15	0.83	0.24	65,69,73,79	0
5	FUC	K	4	10/11	0.84	0.30	59,64,68,68	0
6	BMA	T	3	11/12	0.84	0.14	51,54,57,60	0
6	NAG	M	1	14/15	0.84	0.20	48,57,61,65	0
8	BMA	S	3	11/12	0.84	0.19	49,51,62,68	0

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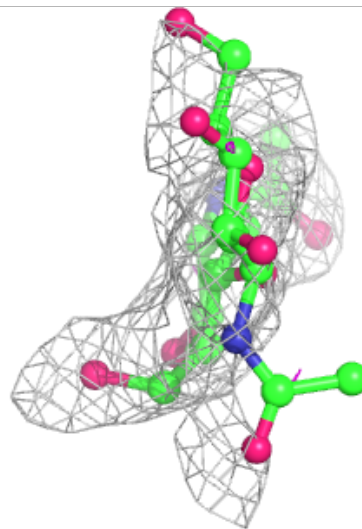
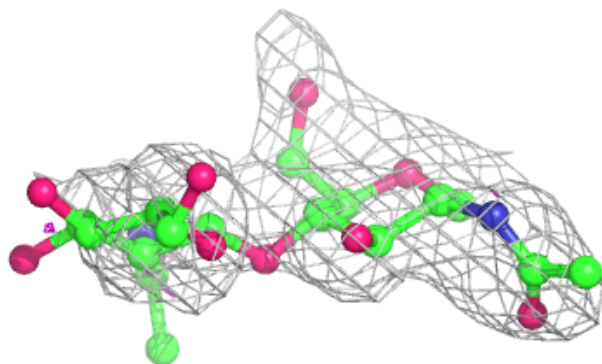
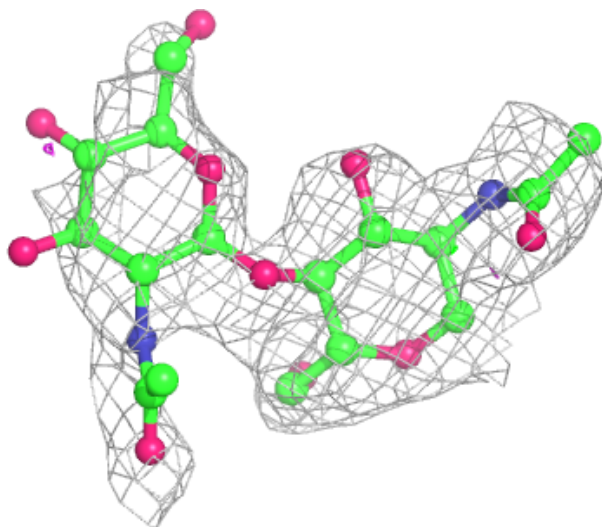
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	Y	1	14/15	0.84	0.23	52,60,68,72	0
3	NAG	a	2	14/15	0.85	0.28	61,67,69,69	0
7	NAG	N	1	14/15	0.85	0.23	54,61,65,68	0
3	NAG	H	1	14/15	0.85	0.20	62,67,74,79	0
9	NAG	U	1	14/15	0.85	0.20	56,64,69,69	0
4	MAN	I	5	11/12	0.86	0.20	82,85,90,91	0
7	FUC	N	2	10/11	0.86	0.35	72,76,79,81	0
3	NAG	G	1	14/15	0.87	0.24	71,75,81,89	0
4	BMA	X	3	11/12	0.87	0.12	61,67,73,75	0
9	NAG	Z	1	14/15	0.87	0.18	56,60,68,72	0
5	NAG	K	1	14/15	0.88	0.15	43,45,55,58	0
8	FUC	S	6	10/11	0.89	0.16	49,52,54,56	0
8	MAN	S	4	11/12	0.90	0.12	44,49,53,54	0
4	NAG	P	1	14/15	0.90	0.24	51,57,60,62	0
3	NAG	a	1	14/15	0.90	0.13	41,48,52,54	0
6	NAG	T	2	14/15	0.90	0.12	38,42,44,48	0
4	NAG	I	2	14/15	0.91	0.11	43,48,54,59	0
4	NAG	X	2	14/15	0.91	0.12	41,47,53,57	0
8	NAG	S	2	14/15	0.91	0.11	39,45,48,49	0
6	NAG	T	1	14/15	0.92	0.13	28,32,35,36	0
4	NAG	Q	1	14/15	0.92	0.12	38,41,42,43	0
4	NAG	Q	2	14/15	0.93	0.11	43,48,52,57	0
4	NAG	X	1	14/15	0.93	0.12	39,42,44,44	0
4	NAG	I	1	14/15	0.93	0.10	37,40,42,43	0
8	NAG	S	1	14/15	0.93	0.09	30,35,41,45	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.

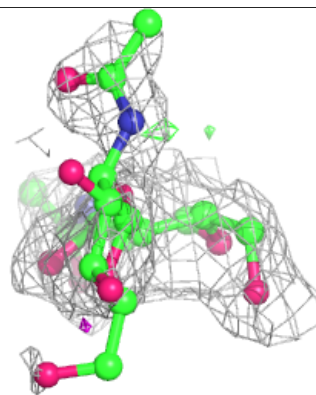
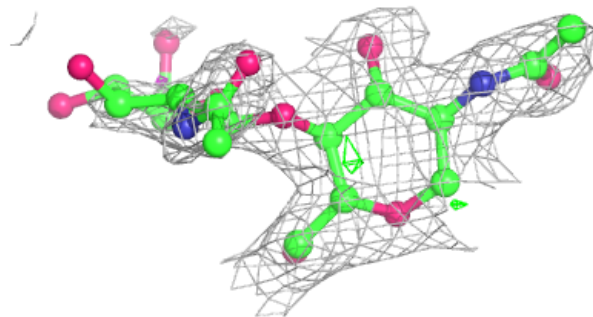
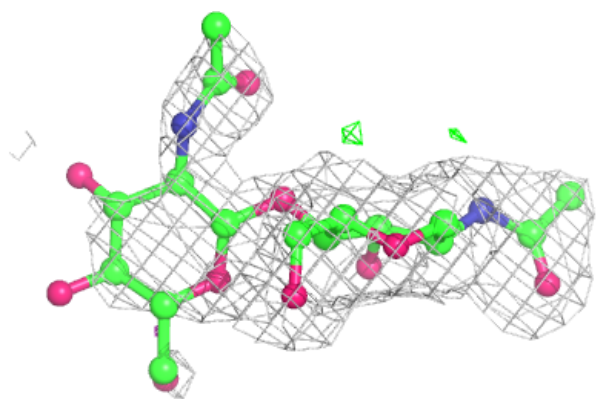
Electron density around Chain G:

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



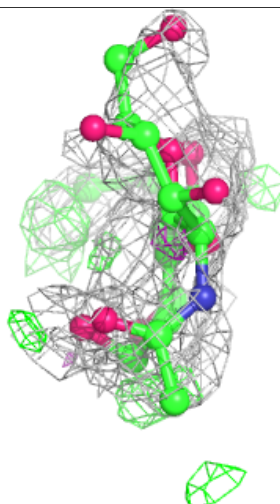
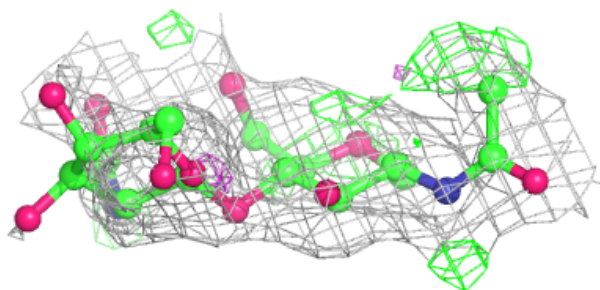
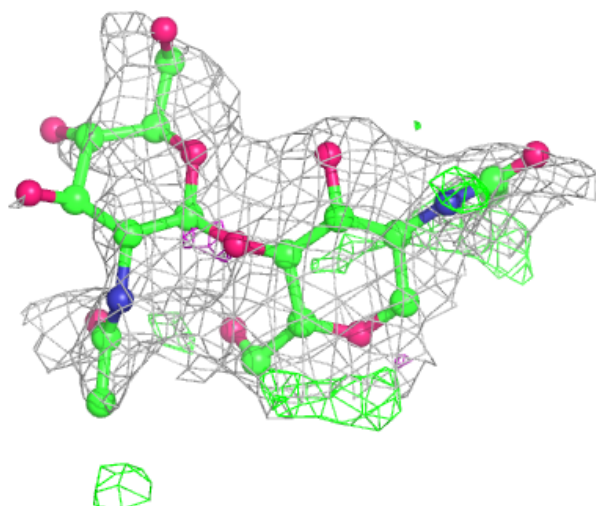
Electron density around Chain H:

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



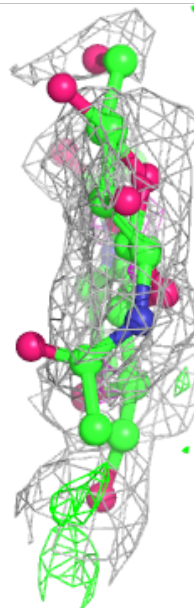
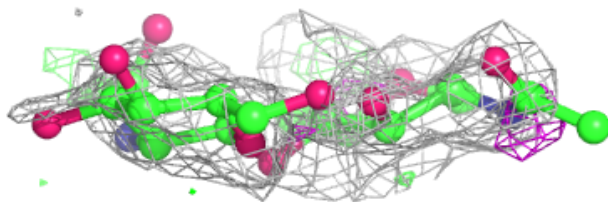
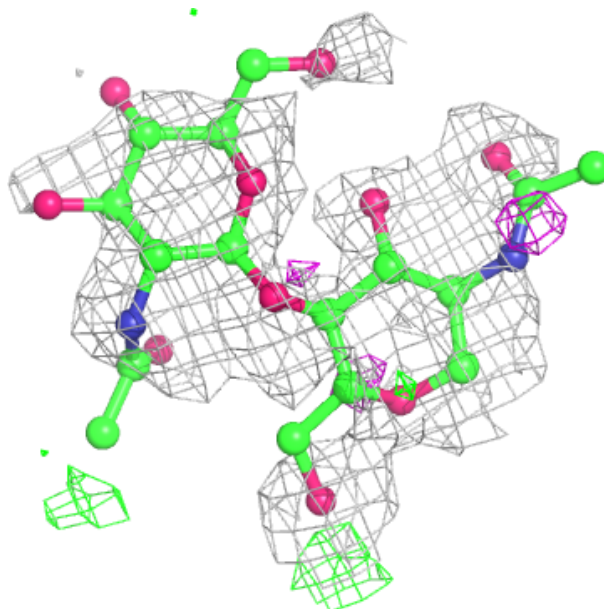
Electron density around Chain J:

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



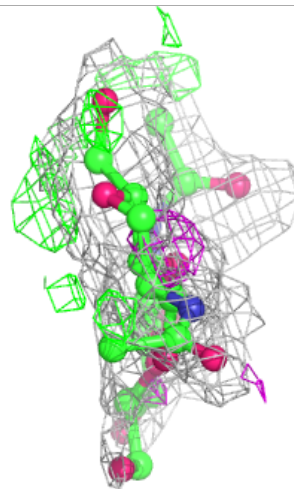
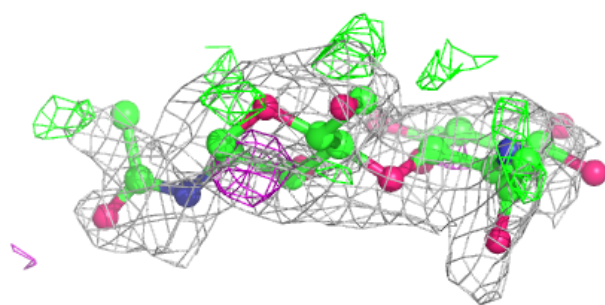
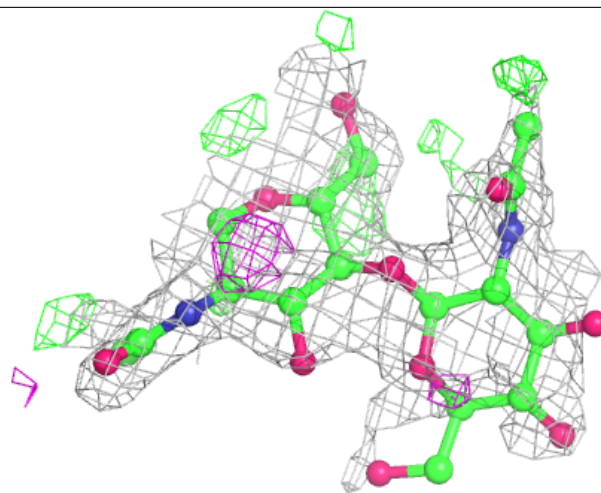
Electron density around Chain L:

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



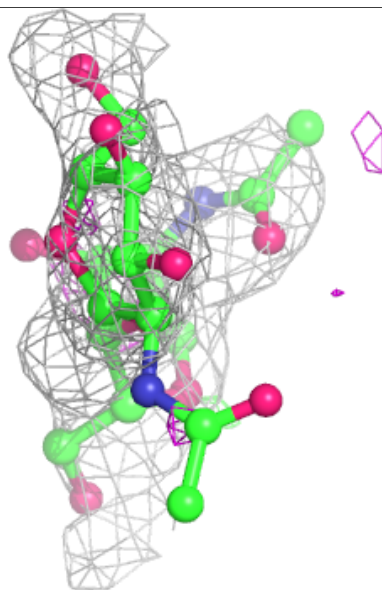
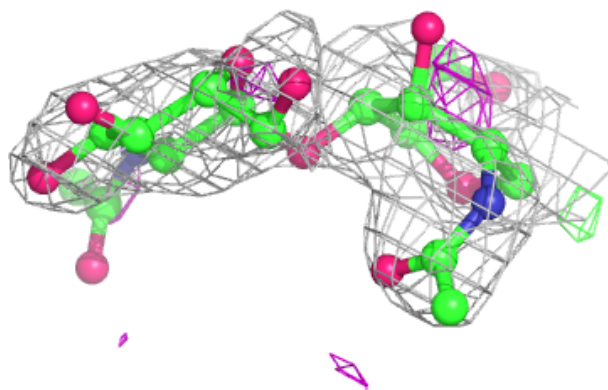
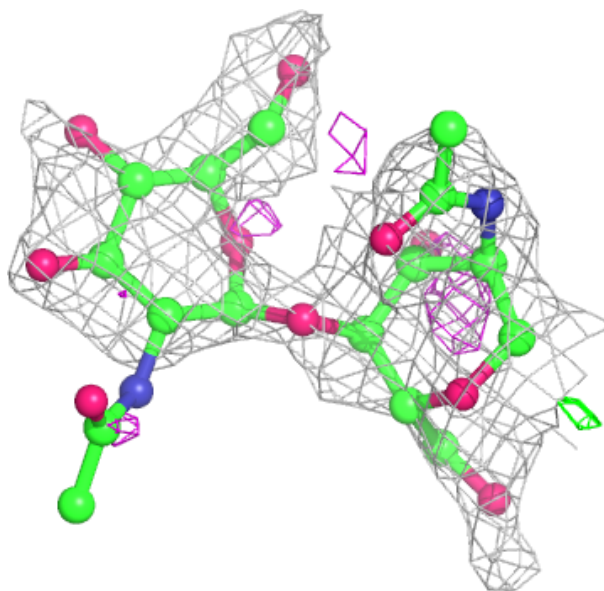
Electron density around Chain R:

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



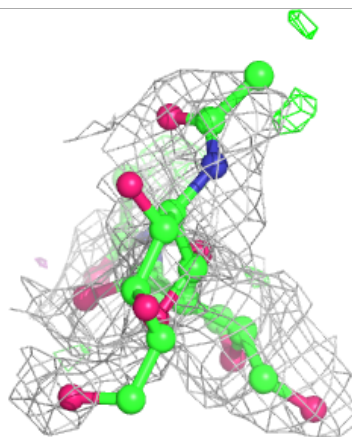
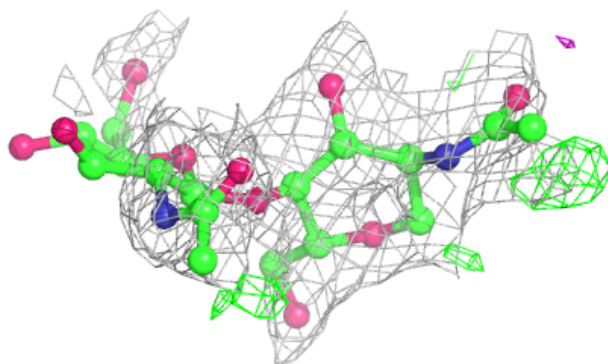
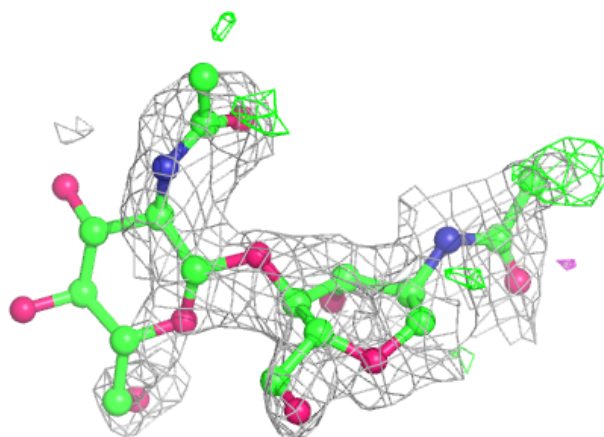
Electron density around Chain V:

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



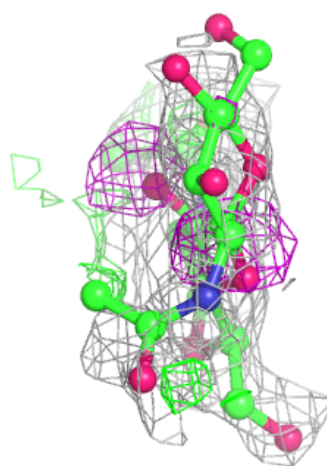
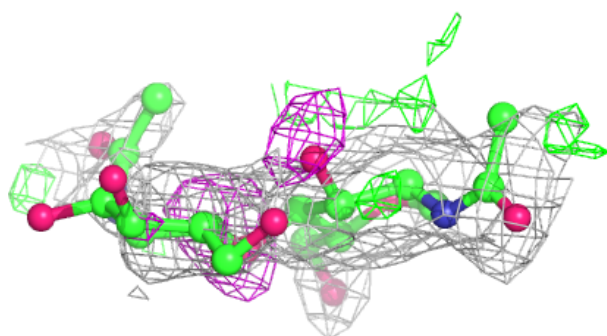
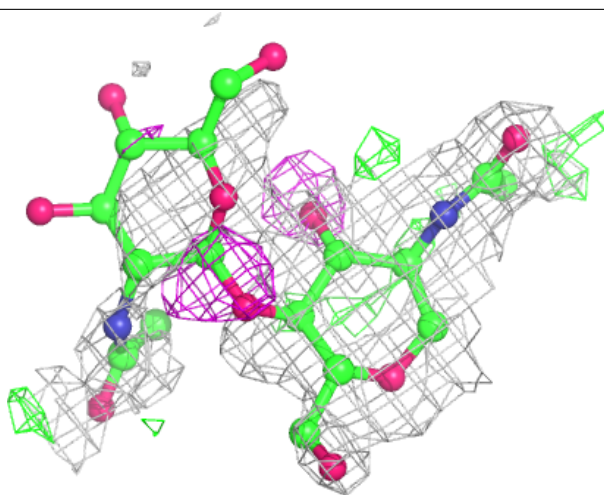
Electron density around Chain W:

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



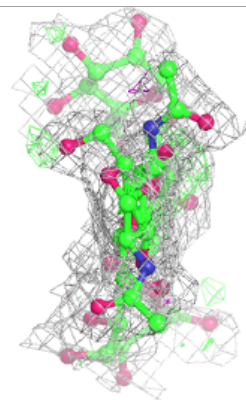
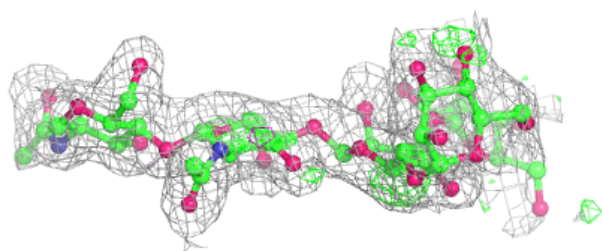
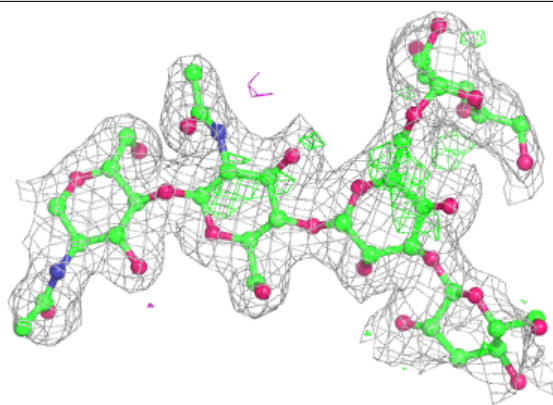
Electron density around Chain Y:

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



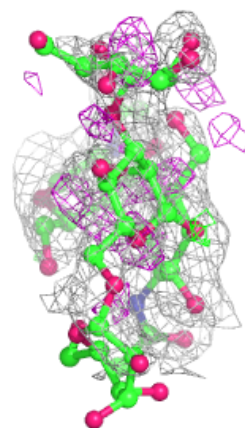
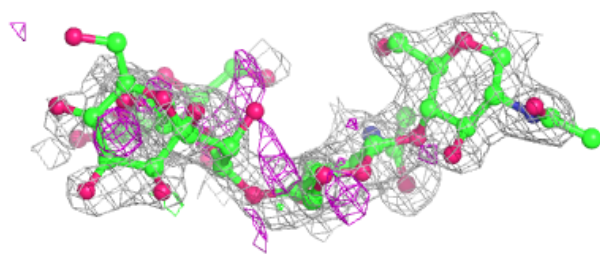
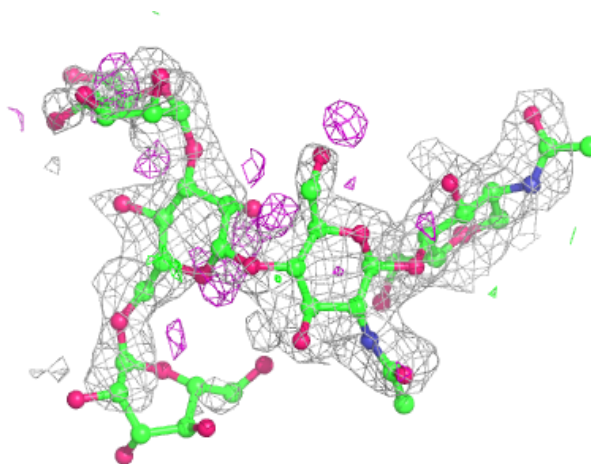
Electron density around Chain I:

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



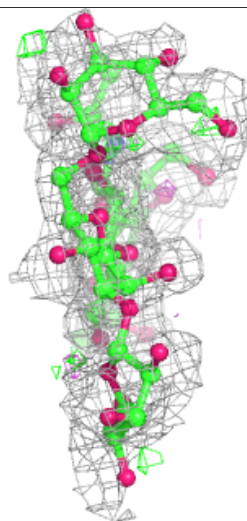
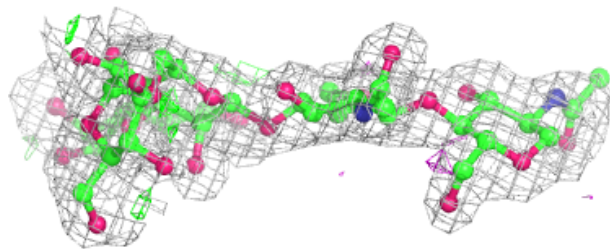
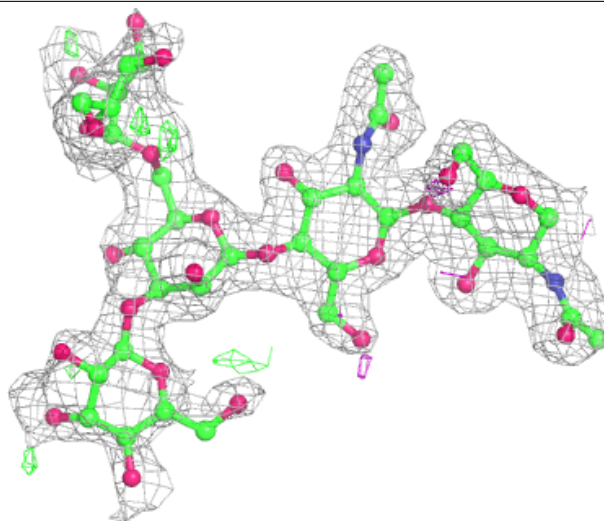
Electron density around Chain P:

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



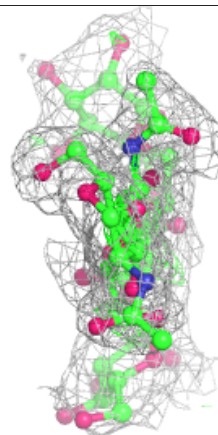
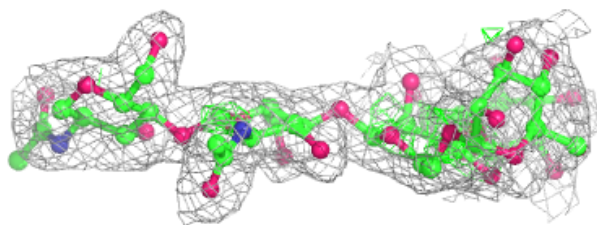
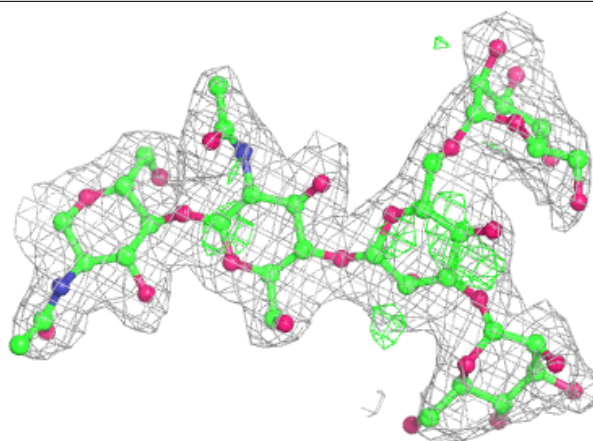
Electron density around Chain Q:

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



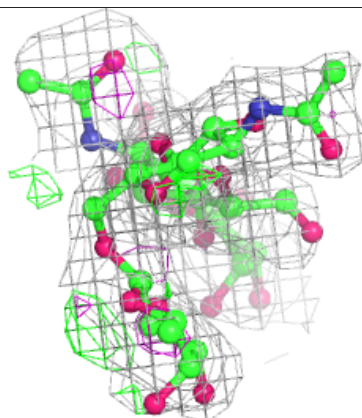
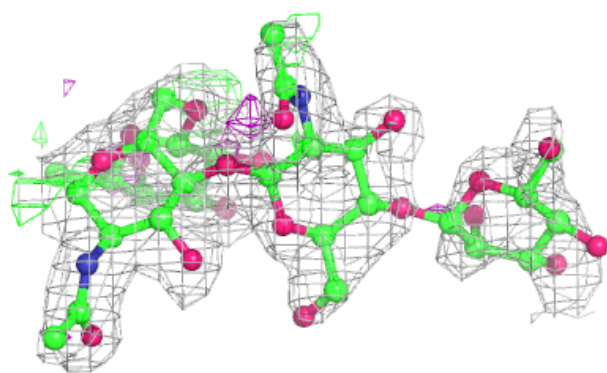
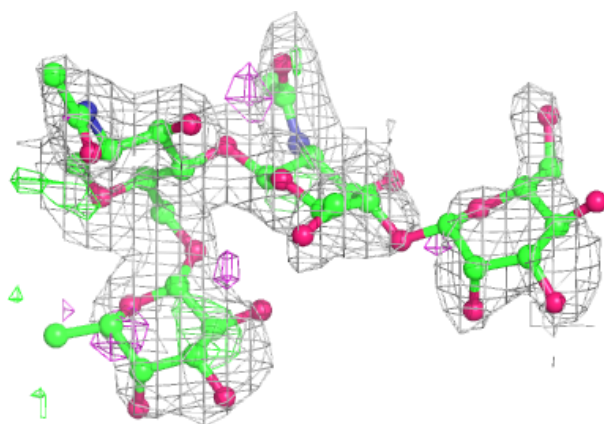
Electron density around Chain X:

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

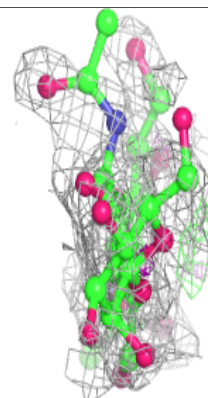
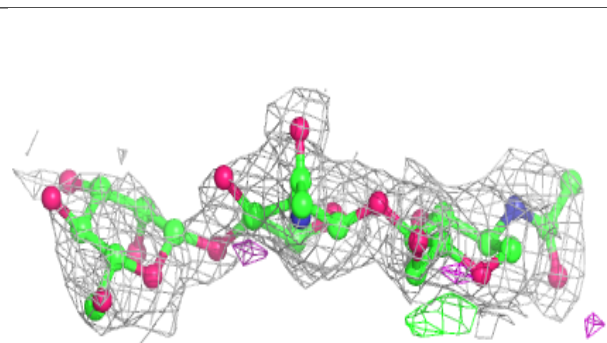
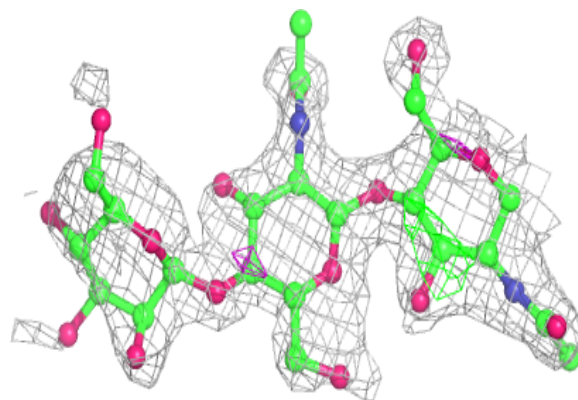


Electron density around Chain K:

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

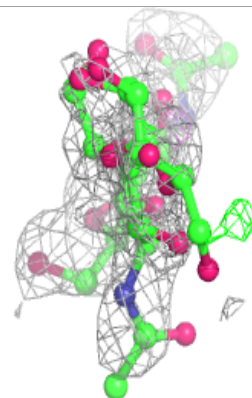
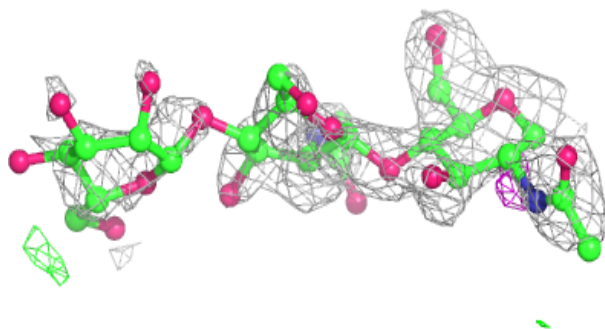
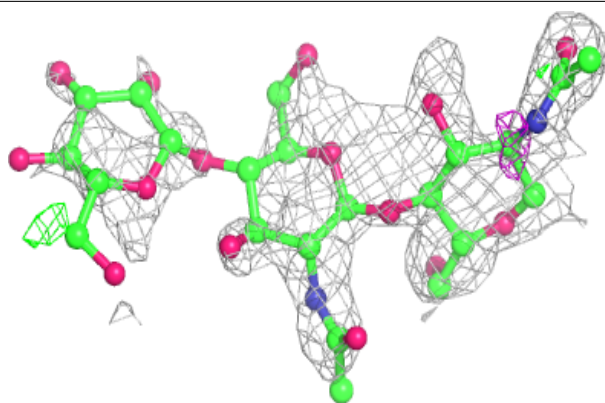
**Electron density around Chain M:**

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

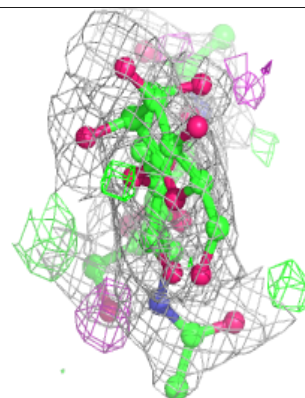
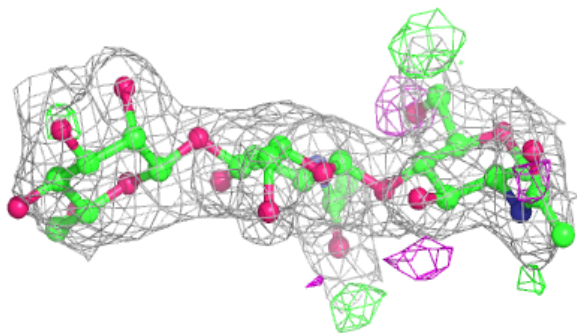
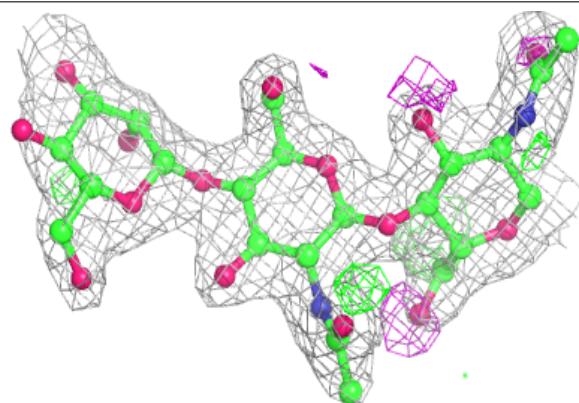


Electron density around Chain O:

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

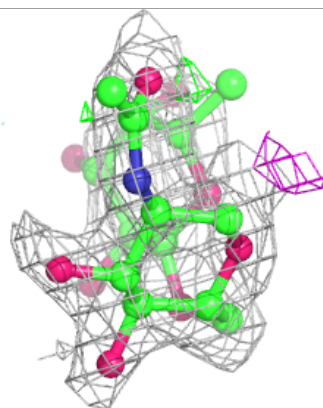
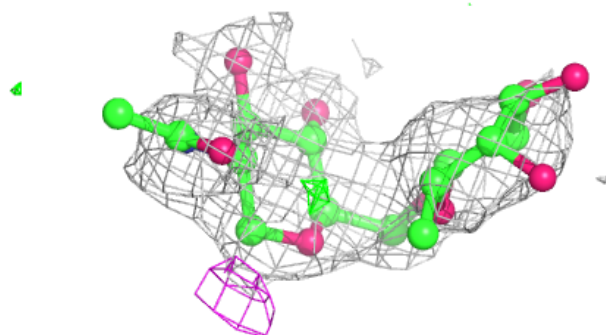
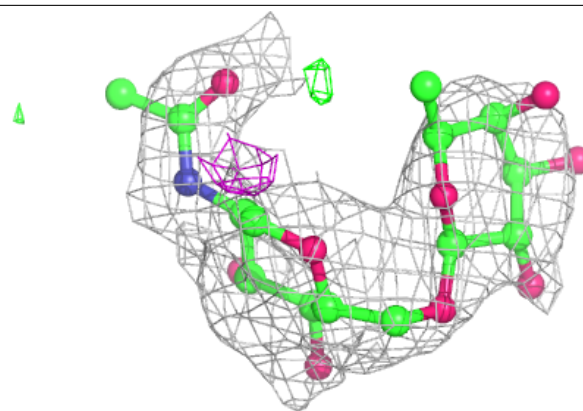
**Electron density around Chain T:**

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

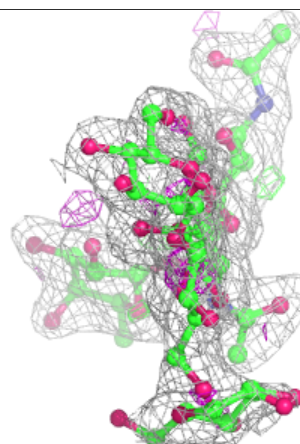
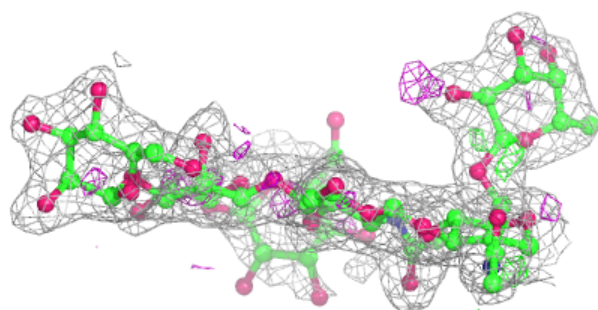
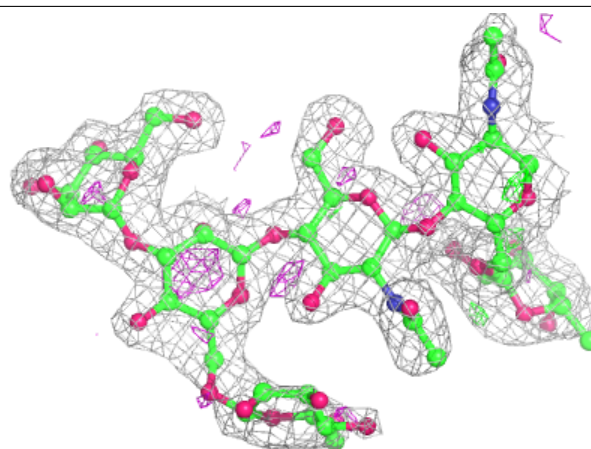


Electron density around Chain N:

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

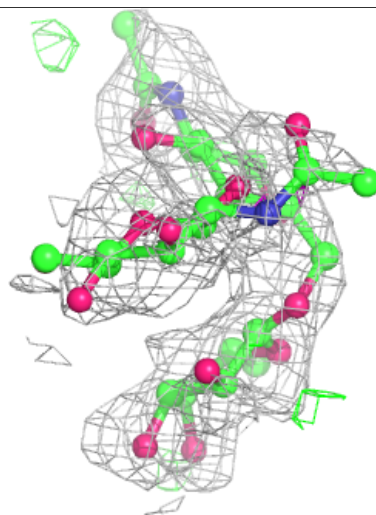
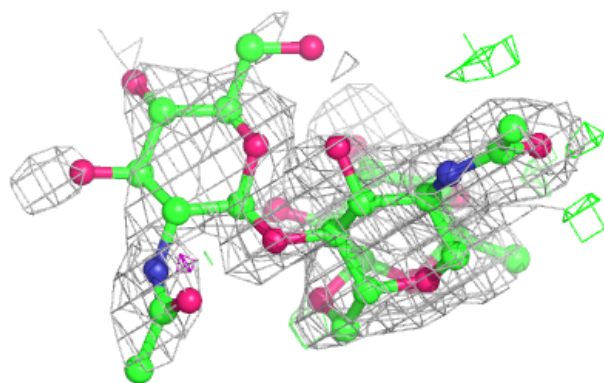
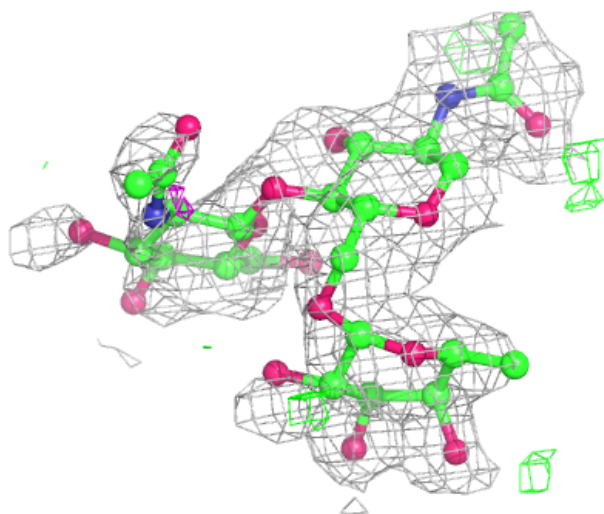
**Electron density around Chain S:**

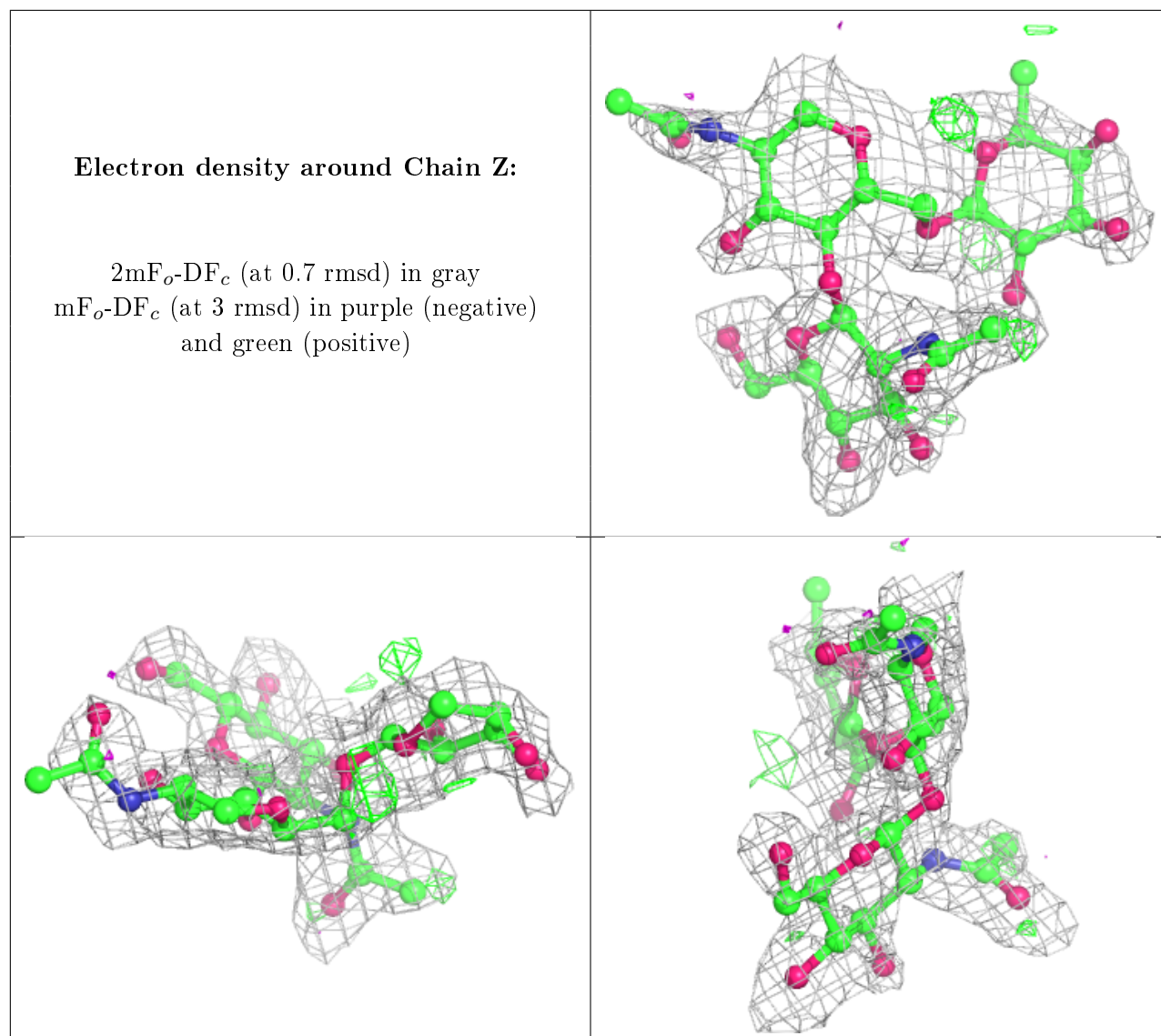
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain U:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
10	BMA	A	604	11/12	0.60	0.52	126,134,137,137	0
12	NAG	C	601	14/15	0.61	0.44	90,93,98,100	0
11	EDO	C	1327	4/4	0.62	0.19	65,65,67,67	0
12	NAG	E	621	14/15	0.71	0.51	68,76,80,81	0
11	EDO	A	1328	4/4	0.77	0.28	61,63,64,65	0
11	EDO	C	1328	4/4	0.81	0.25	35,38,42,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
11	EDO	E	1327	4/4	0.82	0.14	48,50,52,53	0
11	EDO	F	1173	4/4	0.85	0.19	58,59,60,61	0
11	EDO	F	1174	4/4	0.86	0.16	53,53,55,56	0
11	EDO	E	1326	4/4	0.86	0.18	54,55,57,61	0
11	EDO	D	1173	4/4	0.88	0.29	51,52,52,53	0
11	EDO	A	1327	4/4	0.88	0.14	37,37,39,40	0
11	EDO	F	1175	4/4	0.88	0.12	48,51,52,52	0
11	EDO	B	1173	4/4	0.90	0.22	40,40,42,45	0
11	EDO	B	1174	4/4	0.93	0.16	44,45,45,45	0

6.5 Other polymers ⓘ

There are no such residues in this entry.