



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 04:34 PM UTC

PDB ID : 5M8M / pdb_00005m8m
Title : Crystal structure of human tyrosinase related protein 1 in complex with kojic acid
Authors : Lai, X.; Soler-Lopez, M.; Wichers, H.J.; Dijkstra, B.W.
Deposited on : 2016-10-29
Resolution : 2.65 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

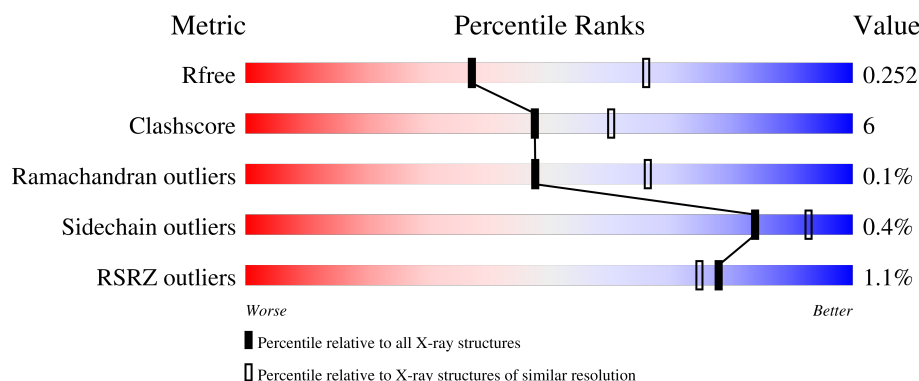
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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.



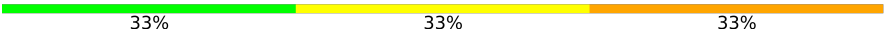
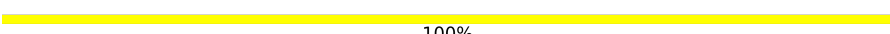
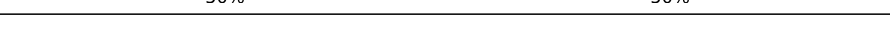
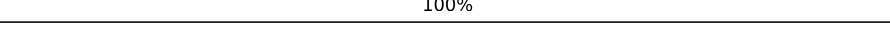
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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

Mol	Chain	Length	Quality of chain
1	A	446	% 89% 11%
1	B	446	2% 87% 12%
1	C	446	% 89% 11%
1	D	446	% 86% 13%
2	E	3	33% 33% 33%

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Mol	Chain	Length	Quality of chain
2	J	3	
2	O	3	
2	T	3	
3	F	2	
3	K	2	
3	U	2	
4	G	2	
4	H	2	
4	I	2	
4	L	2	
4	N	2	
4	Q	2	
4	S	2	
4	V	2	
4	W	2	
4	X	2	
5	M	4	
6	P	5	
7	R	5	

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
9	KOJ	A	503	-	X	-	-
9	KOJ	D	504	-	X	-	-

2 Entry composition [i](#)

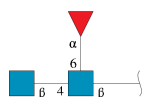
There are 11 unique types of molecules in this entry. The entry contains 15021 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 5,6-dihydroxyindole-2-carboxylic acid oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	446	Total	C	N	O	S	0	0	0
			3560	2233	632	672	23			
1	B	446	Total	C	N	O	S	0	0	0
			3560	2233	632	672	23			
1	C	446	Total	C	N	O	S	0	0	0
			3560	2233	632	672	23			
1	D	446	Total	C	N	O	S	0	0	0
			3560	2233	632	672	23			

- Molecule 2 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
2	E	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	J	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	O	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	T	3	Total	C	N	O	0	0	0
			38	22	2	14			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	2	Total	C	N	O	0	0	0
			24	14	1	9			
3	K	2	Total	C	N	O	0	0	0
			24	14	1	9			
3	U	2	Total	C	N	O	0	0	0
			24	14	1	9			

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



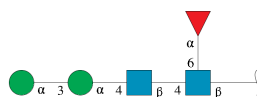
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	L	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	N	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	Q	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	S	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	V	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	W	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	X	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



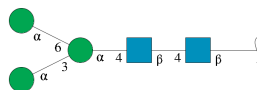
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	M	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-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
6	P	5	Total	C	N	O	0	0	0
			60	34	2	24			

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

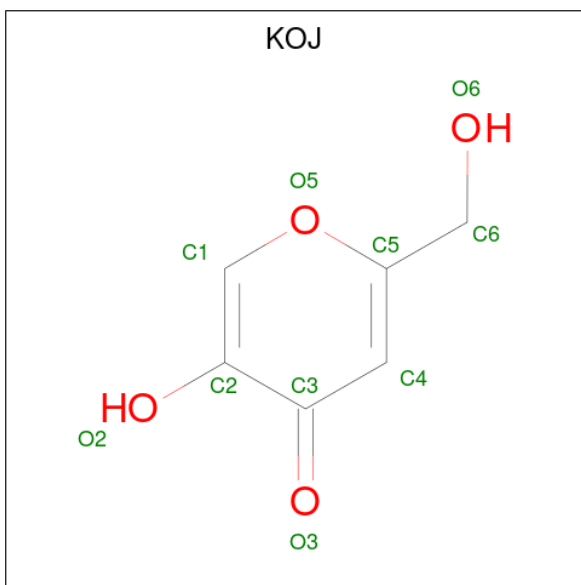


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	R	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 8 is ZINC ION (CCD ID: ZN) (formula: Zn).

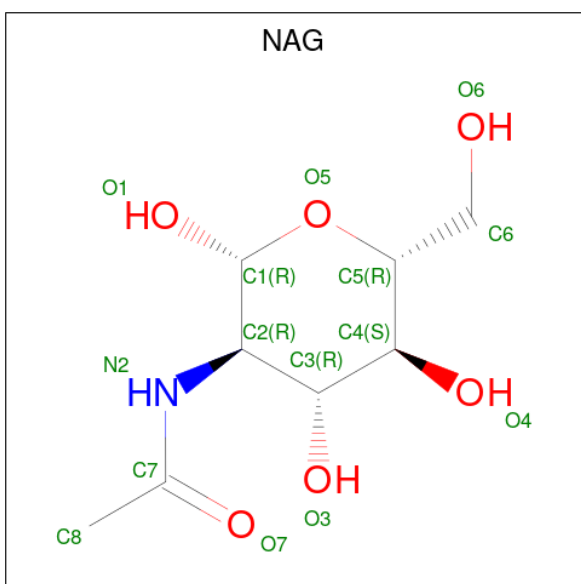
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	2	Total	Zn	0	0
			2	2		
8	B	3	Total	Zn	0	0
			3	3		
8	C	2	Total	Zn	0	0
			2	2		
8	D	2	Total	Zn	0	0
			2	2		

- Molecule 9 is 5-HYDROXY-2-(HYDROXYMETHYL)-4H-PYRAN-4-ONE (CCD ID: KOJ) (formula: C₆H₆O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			10	6	4		
9	B	1	Total	C	O	0	0
			10	6	4		
9	C	1	Total	C	O	0	0
			10	6	4		
9	D	1	Total	C	O	0	0
			10	6	4		

- Molecule 10 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

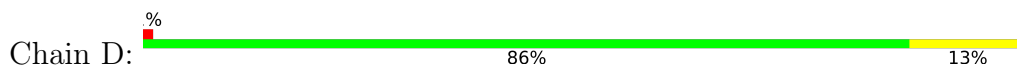


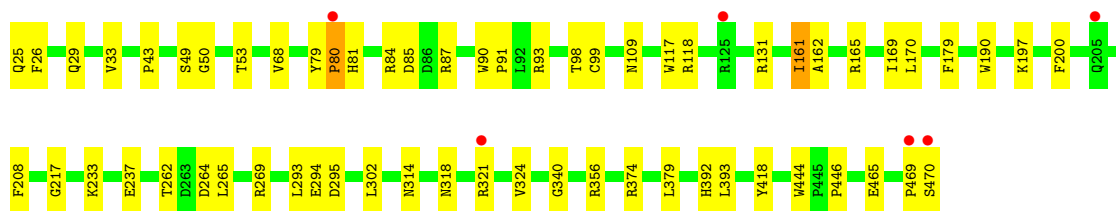
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	17	Total	O	0	0
			17	17		
11	B	7	Total	O	0	0
			7	7		
11	C	12	Total	O	0	0
			12	12		
11	D	7	Total	O	0	0
			7	7		

- Molecule 1: 5,6-dihydroxyindole-2-carboxylic acid oxidase





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

Chain E: 33% 33% 33%



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

Chain J: 33% 33% 33%



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

Chain O: 33% 67%



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

Chain T: 33% 67%



- Molecule 3: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F: 100%



- Molecule 3: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K: 100%

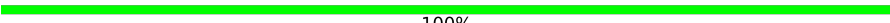


- Molecule 3: α -L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain U:  100%

NA G1
FUC2

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

Chain G:  100%

NA G1
NA G2

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

Chain H:  50% 50%

NA G1
NA G2

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

Chain I:  100%

NA G1
NA G2

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

Chain L:  50% 50%

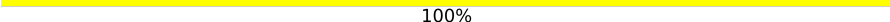
NA G1
NA G2

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

Chain N:  50% 50%

NA G1
NA G2

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

Chain Q:  100%

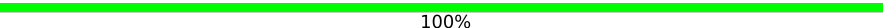
MAG1
MAG2

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

Chain S:  50% 50%

MAG1
MAG2

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

Chain V:  100%

MAG1
MAG2

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

Chain W:  50% 50%

MAG1
MAG2

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

Chain X:  50% 50%

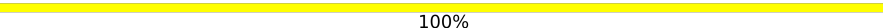
MAG1
MAG2

- Molecule 5: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  50% 50%

MAG1
MAG2
MAN3
MAN4

- Molecule 6: alpha-D-mannopyranose-(1-3)-alpha-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 P:  100%

MAG1
MAG2
MAN3
MAN4
FUC5

- Molecule 7: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] α -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain R:



MAG1
MAG2
MAG3
MAG4
MAG5

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.25Å 141.53Å 192.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.00 – 2.65 49.00 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.00-2.65) 93.0 (49.00-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.65Å)	Xtriage
Refinement program	REFMAC, PHENIX	Depositor
R, R_{free}	0.197 , 0.249 0.205 , 0.252	Depositor DCC
R_{free} test set	3534 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	42.7	Xtriage
Anisotropy	0.493	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15021	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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: KOJ, MAN, NAG, FUC, ZN

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	0.55	0/3667	0.58	1/4998 (0.0%)
1	B	0.51	1/3667 (0.0%)	0.59	3/4998 (0.1%)
1	C	0.56	1/3667 (0.0%)	0.59	0/4998
1	D	0.51	1/3667 (0.0%)	0.56	1/4998 (0.0%)
All	All	0.53	3/14668 (0.0%)	0.58	5/19992 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	80	PRO	N-CA	7.11	1.56	1.47
1	B	381	HIS	C-O	-5.54	1.17	1.24
1	C	381	HIS	C-O	-5.36	1.17	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	118	ARG	NE-CZ-NH2	-10.32	109.91	119.20
1	B	126	VAL	CA-CB-CG2	7.39	122.96	110.40
1	B	118	ARG	NH1-CZ-NH2	5.95	127.03	119.30
1	D	80	PRO	N-CA-CB	-5.54	97.43	103.25
1	A	345	PRO	O-C-N	5.07	123.53	121.15

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	3560	0	3326	35	0
1	B	3560	0	3326	37	0
1	C	3560	0	3326	37	0
1	D	3560	0	3323	47	0
2	E	38	0	34	1	0
2	J	38	0	34	1	0
2	O	38	0	34	2	0
2	T	38	0	33	0	0
3	F	24	0	22	0	0
3	K	24	0	22	0	0
3	U	24	0	22	1	0
4	G	28	0	25	0	0
4	H	28	0	25	1	0
4	I	28	0	25	0	0
4	L	28	0	25	0	0
4	N	28	0	25	2	0
4	Q	28	0	25	0	0
4	S	28	0	25	0	0
4	V	28	0	25	0	0
4	W	28	0	25	2	0
4	X	28	0	25	1	0
5	M	50	0	43	1	0
6	P	60	0	52	0	0
7	R	61	0	52	3	0
8	A	2	0	0	0	0
8	B	3	0	0	0	0
8	C	2	0	0	0	0
8	D	2	0	0	0	0
9	A	10	0	6	0	0
9	B	10	0	5	0	0
9	C	10	0	5	0	0
9	D	10	0	5	1	0
10	D	14	0	13	0	0
11	A	17	0	0	0	0
11	B	7	0	0	0	0
11	C	12	0	0	1	0
11	D	7	0	0	1	0
All	All	15021	0	13933	162	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 6.

The worst 5 of 162 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:C:118:ARG:NE	1:C:126:VAL:HG11	1.89	0.86
1:D:80:PRO:HG2	1:D:81:HIS:ND1	1.96	0.79
1:B:203:VAL:HB	1:C:241:GLU:HG2	1.66	0.77
1:B:64:ARG:NH2	1:B:66:GLU:OE2	2.17	0.76
7:R:2:NAG:H3	7:R:2:NAG:H82	1.68	0.75

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	444/446 (100%)	421 (95%)	23 (5%)	0	100	100
1	B	444/446 (100%)	420 (95%)	24 (5%)	0	100	100
1	C	444/446 (100%)	422 (95%)	21 (5%)	1 (0%)	43	60
1	D	444/446 (100%)	420 (95%)	24 (5%)	0	100	100
All	All	1776/1784 (100%)	1683 (95%)	92 (5%)	1 (0%)	48	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	376	LEU

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	395/395 (100%)	391 (99%)	4 (1%)	68	81
1	B	395/395 (100%)	393 (100%)	2 (0%)	81	91
1	C	395/395 (100%)	395 (100%)	0	100	100
1	D	395/395 (100%)	394 (100%)	1 (0%)	86	94
All	All	1580/1580 (100%)	1573 (100%)	7 (0%)	84	93

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	349	SER
1	B	349	SER
1	D	161	ILE
1	B	447	VAL
1	A	168	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	78	GLN
1	D	325	GLN
1	D	353	ASN
1	C	104	ASN
1	C	142	ASN

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 ⓘ

52 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	E	1	1,2	14,14,15	0.98	1 (7%)	17,19,21	0.64	0
2	NAG	E	2	2	14,14,15	0.58	0	17,19,21	0.56	0
2	FUC	E	3	2	10,10,11	0.93	0	14,14,16	1.42	3 (21%)
3	NAG	F	1	1,3	14,14,15	1.51	2 (14%)	17,19,21	0.99	1 (5%)
3	FUC	F	2	3	10,10,11	1.38	2 (20%)	14,14,16	1.42	3 (21%)
4	NAG	G	1	1,4	14,14,15	0.36	0	17,19,21	0.63	0
4	NAG	G	2	4	14,14,15	0.24	0	17,19,21	0.36	0
4	NAG	H	1	1,4	14,14,15	0.52	0	17,19,21	0.81	1 (5%)
4	NAG	H	2	4	14,14,15	1.66	2 (14%)	17,19,21	1.52	3 (17%)
4	NAG	I	1	1,4	14,14,15	0.49	0	17,19,21	0.55	0
4	NAG	I	2	4	14,14,15	0.38	0	17,19,21	0.39	0
2	NAG	J	1	1,2	14,14,15	0.58	1 (7%)	17,19,21	0.60	0
2	NAG	J	2	2	14,14,15	0.55	0	17,19,21	0.47	0
2	FUC	J	3	2	10,10,11	0.66	0	14,14,16	1.34	2 (14%)
3	NAG	K	1	1,3	14,14,15	1.73	3 (21%)	17,19,21	1.17	2 (11%)
3	FUC	K	2	3	10,10,11	1.20	2 (20%)	14,14,16	1.26	3 (21%)
4	NAG	L	1	1,4	14,14,15	0.60	0	17,19,21	0.76	1 (5%)
4	NAG	L	2	4	14,14,15	0.37	0	17,19,21	0.54	0
5	NAG	M	1	1,5	14,14,15	0.42	0	17,19,21	0.88	1 (5%)
5	NAG	M	2	5	14,14,15	0.92	1 (7%)	17,19,21	1.51	4 (23%)
5	MAN	M	3	5	11,11,12	1.98	3 (27%)	15,15,17	1.94	4 (26%)
5	MAN	M	4	5	11,11,12	1.30	2 (18%)	15,15,17	1.35	4 (26%)
4	NAG	N	1	1,4	14,14,15	0.54	0	17,19,21	0.77	1 (5%)
4	NAG	N	2	4	14,14,15	0.56	0	17,19,21	0.45	0
2	NAG	O	1	1,2	14,14,15	0.62	1 (7%)	17,19,21	0.66	0
2	NAG	O	2	2	14,14,15	0.46	0	17,19,21	0.43	0
2	FUC	O	3	2	10,10,11	0.71	0	14,14,16	1.41	2 (14%)
6	NAG	P	1	1,6	14,14,15	0.71	1 (7%)	17,19,21	0.74	0
6	NAG	P	2	6	14,14,15	1.22	1 (7%)	17,19,21	0.81	0
6	MAN	P	3	6	11,11,12	1.03	2 (18%)	15,15,17	1.67	3 (20%)
6	MAN	P	4	6	11,11,12	1.13	1 (9%)	15,15,17	2.47	4 (26%)
6	FUC	P	5	6	10,10,11	1.51	1 (10%)	14,14,16	1.73	3 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	Q	1	1,4	14,14,15	0.21	0	17,19,21	0.89	1 (5%)
4	NAG	Q	2	4	14,14,15	1.08	2 (14%)	17,19,21	0.81	1 (5%)
7	NAG	R	1	1,7	14,14,15	0.36	0	17,19,21	0.92	2 (11%)
7	NAG	R	2	7	14,14,15	1.21	2 (14%)	17,19,21	1.89	3 (17%)
7	MAN	R	3	7	11,11,12	1.52	2 (18%)	15,15,17	1.88	4 (26%)
7	MAN	R	4	7	11,11,12	1.19	1 (9%)	15,15,17	1.79	3 (20%)
7	MAN	R	5	7	11,11,12	1.10	1 (9%)	15,15,17	1.64	3 (20%)
4	NAG	S	1	1,4	14,14,15	0.77	1 (7%)	17,19,21	0.47	0
4	NAG	S	2	4	14,14,15	0.58	0	17,19,21	0.48	0
2	NAG	T	1	1,2	14,14,15	0.59	1 (7%)	17,19,21	0.46	0
2	NAG	T	2	2	14,14,15	0.64	0	17,19,21	0.46	0
2	FUC	T	3	2	10,10,11	1.14	0	14,14,16	1.33	2 (14%)
3	NAG	U	1	1,3	14,14,15	1.26	2 (14%)	17,19,21	0.99	2 (11%)
3	FUC	U	2	3	10,10,11	1.59	2 (20%)	14,14,16	1.58	3 (21%)
4	NAG	V	1	1,4	14,14,15	0.47	0	17,19,21	0.45	0
4	NAG	V	2	4	14,14,15	0.26	0	17,19,21	0.78	0
4	NAG	W	1	1,4	14,14,15	0.29	0	17,19,21	0.69	0
4	NAG	W	2	4	14,14,15	1.48	2 (14%)	17,19,21	1.56	3 (17%)
4	NAG	X	1	1,4	14,14,15	0.37	0	17,19,21	0.52	0
4	NAG	X	2	4	14,14,15	0.29	0	17,19,21	0.45	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
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	FUC	E	3	2	-	-	0/1/1/1
3	NAG	F	1	1,3	-	2/6/23/26	0/1/1/1
3	FUC	F	2	3	-	-	0/1/1/1
4	NAG	G	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
4	NAG	H	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1
4	NAG	I	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	J	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	FUC	J	3	2	-	-	0/1/1/1
3	NAG	K	1	1,3	-	3/6/23/26	0/1/1/1
3	FUC	K	2	3	-	-	0/1/1/1
4	NAG	L	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	L	2	4	-	2/6/23/26	0/1/1/1
5	NAG	M	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	M	2	5	-	2/6/23/26	0/1/1/1
5	MAN	M	3	5	-	2/2/19/22	1/1/1/1
5	MAN	M	4	5	-	2/2/19/22	1/1/1/1
4	NAG	N	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	N	2	4	-	2/6/23/26	0/1/1/1
2	NAG	O	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	O	2	2	-	2/6/23/26	0/1/1/1
2	FUC	O	3	2	-	-	0/1/1/1
6	NAG	P	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	P	2	6	-	2/6/23/26	0/1/1/1
6	MAN	P	3	6	-	2/2/19/22	1/1/1/1
6	MAN	P	4	6	-	1/2/19/22	0/1/1/1
6	FUC	P	5	6	-	-	0/1/1/1
4	NAG	Q	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	Q	2	4	-	2/6/23/26	0/1/1/1
7	NAG	R	1	1,7	-	4/6/23/26	0/1/1/1
7	NAG	R	2	7	-	5/6/23/26	0/1/1/1
7	MAN	R	3	7	-	2/2/19/22	1/1/1/1
7	MAN	R	4	7	-	2/2/19/22	0/1/1/1
7	MAN	R	5	7	-	1/2/19/22	0/1/1/1
4	NAG	S	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	S	2	4	-	0/6/23/26	0/1/1/1
2	NAG	T	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	T	2	2	-	0/6/23/26	0/1/1/1
2	FUC	T	3	2	-	-	0/1/1/1
3	NAG	U	1	1,3	-	4/6/23/26	0/1/1/1
3	FUC	U	2	3	-	-	0/1/1/1
4	NAG	V	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	V	2	4	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	W	1	1,4	-	4/6/23/26	0/1/1/1
4	NAG	W	2	4	-	3/6/23/26	0/1/1/1
4	NAG	X	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	X	2	4	-	2/6/23/26	0/1/1/1

The worst 5 of 42 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	K	1	NAG	O5-C1	-5.30	1.34	1.43
4	H	2	NAG	C1-C2	5.06	1.59	1.52
4	W	2	NAG	C1-C2	4.61	1.58	1.52
3	F	1	NAG	O5-C1	-4.59	1.36	1.43
5	M	3	MAN	C2-C3	-4.52	1.45	1.52

The worst 5 of 72 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	P	4	MAN	C1-O5-C5	7.14	121.75	112.19
7	R	2	NAG	C2-N2-C7	5.13	129.77	122.90
7	R	4	MAN	C1-O5-C5	4.62	118.38	112.19
7	R	3	MAN	O2-C2-C3	-4.60	100.62	110.15
5	M	3	MAN	C1-O5-C5	4.44	118.13	112.19

There are no chirality outliers.

5 of 79 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	J	2	NAG	O5-C5-C6-O6
4	W	1	NAG	O5-C5-C6-O6
6	P	3	MAN	C4-C5-C6-O6
5	M	4	MAN	O5-C5-C6-O6
4	V	2	NAG	O5-C5-C6-O6

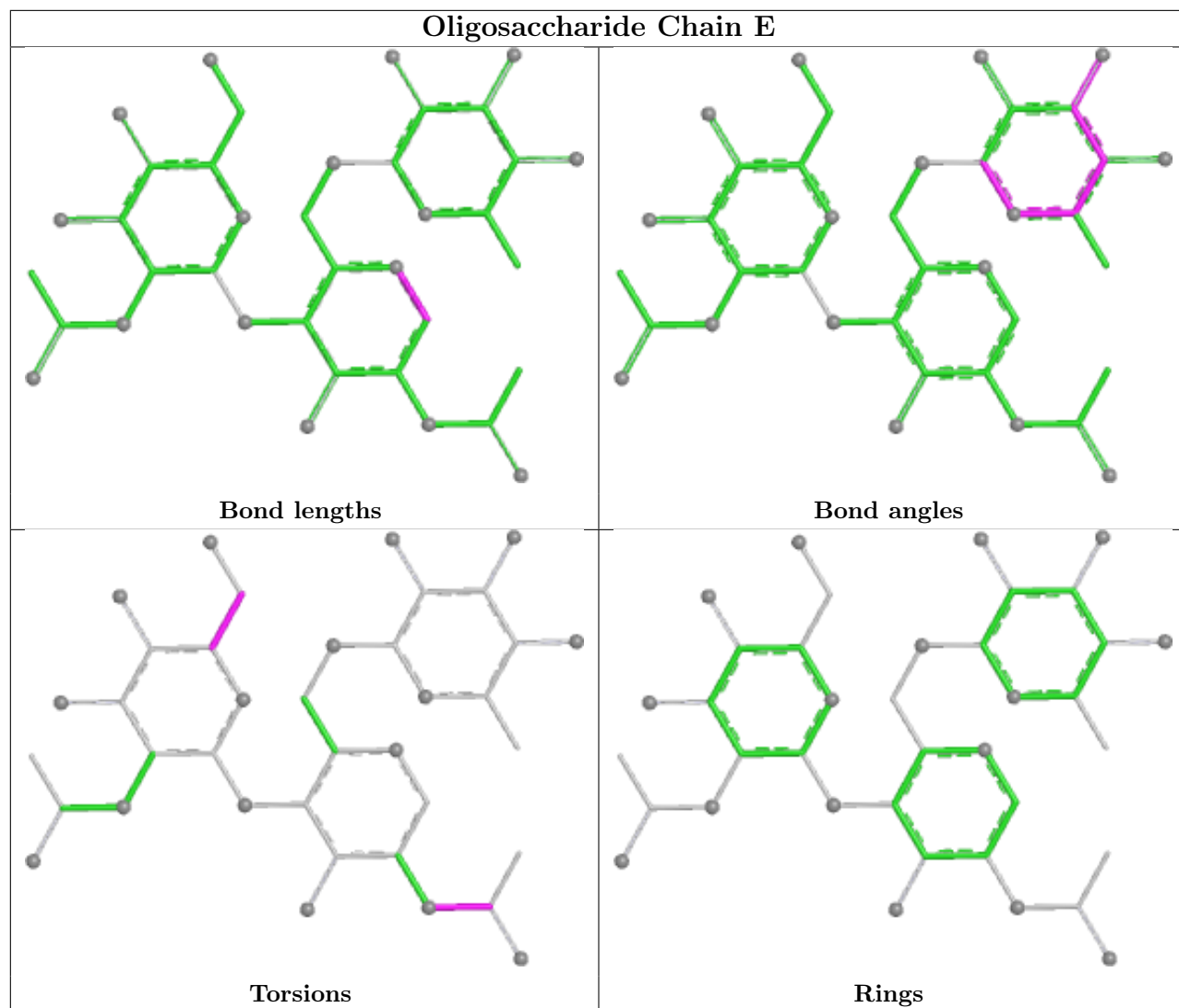
All (4) ring outliers are listed below:

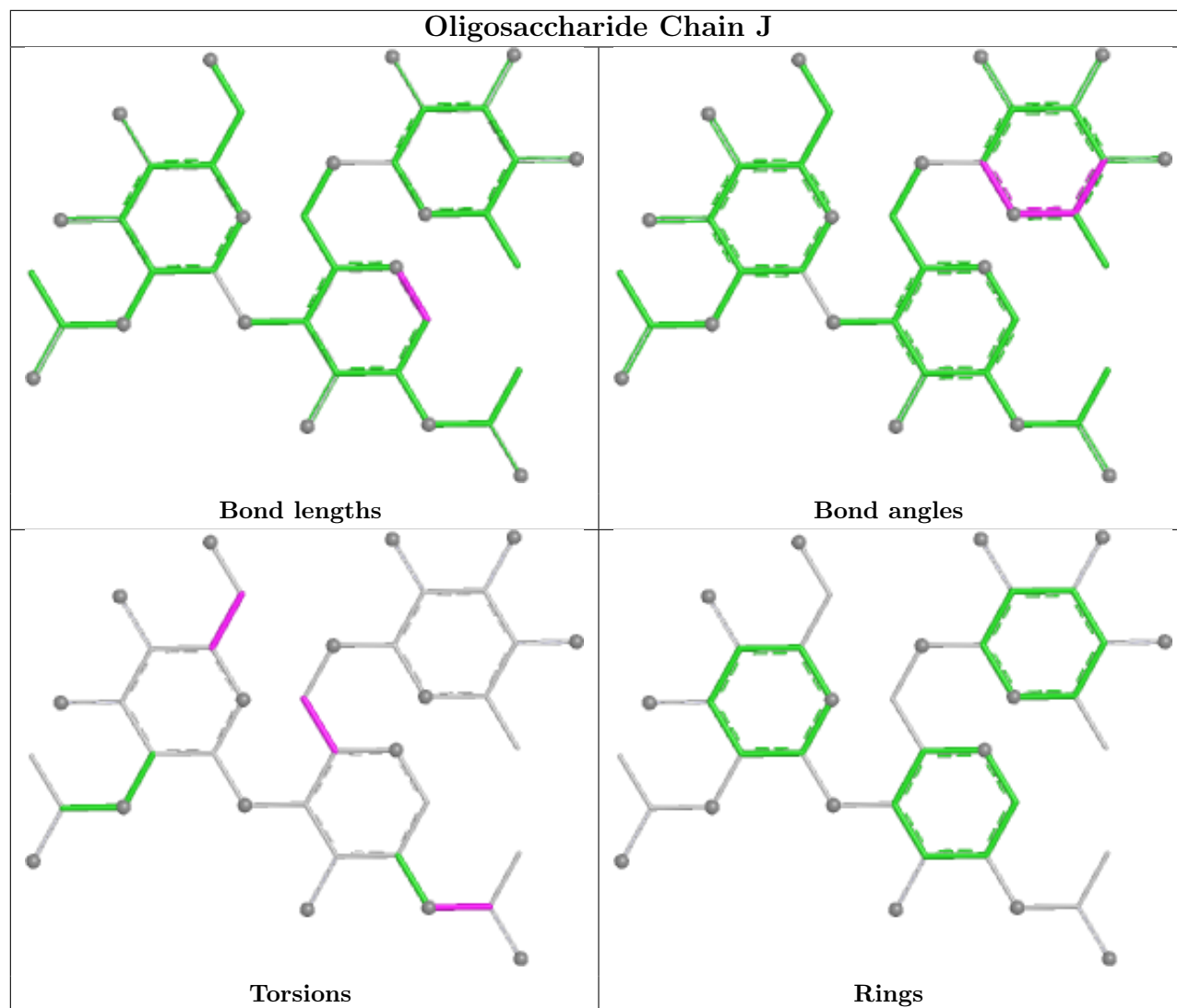
Mol	Chain	Res	Type	Atoms
5	M	3	MAN	C1-C2-C3-C4-C5-O5
6	P	3	MAN	C1-C2-C3-C4-C5-O5
7	R	3	MAN	C1-C2-C3-C4-C5-O5
5	M	4	MAN	C1-C2-C3-C4-C5-O5

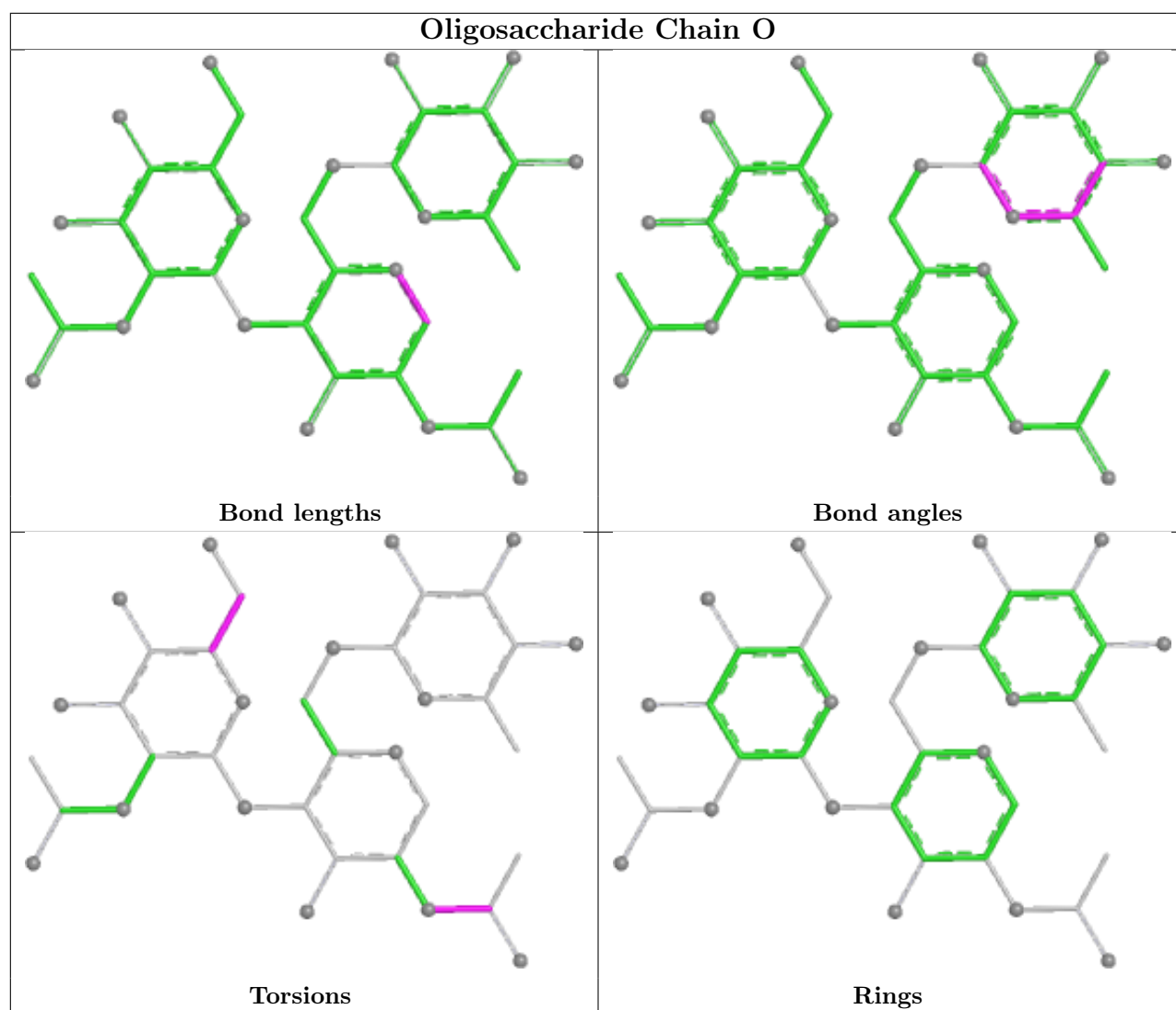
15 monomers are involved in 15 short contacts:

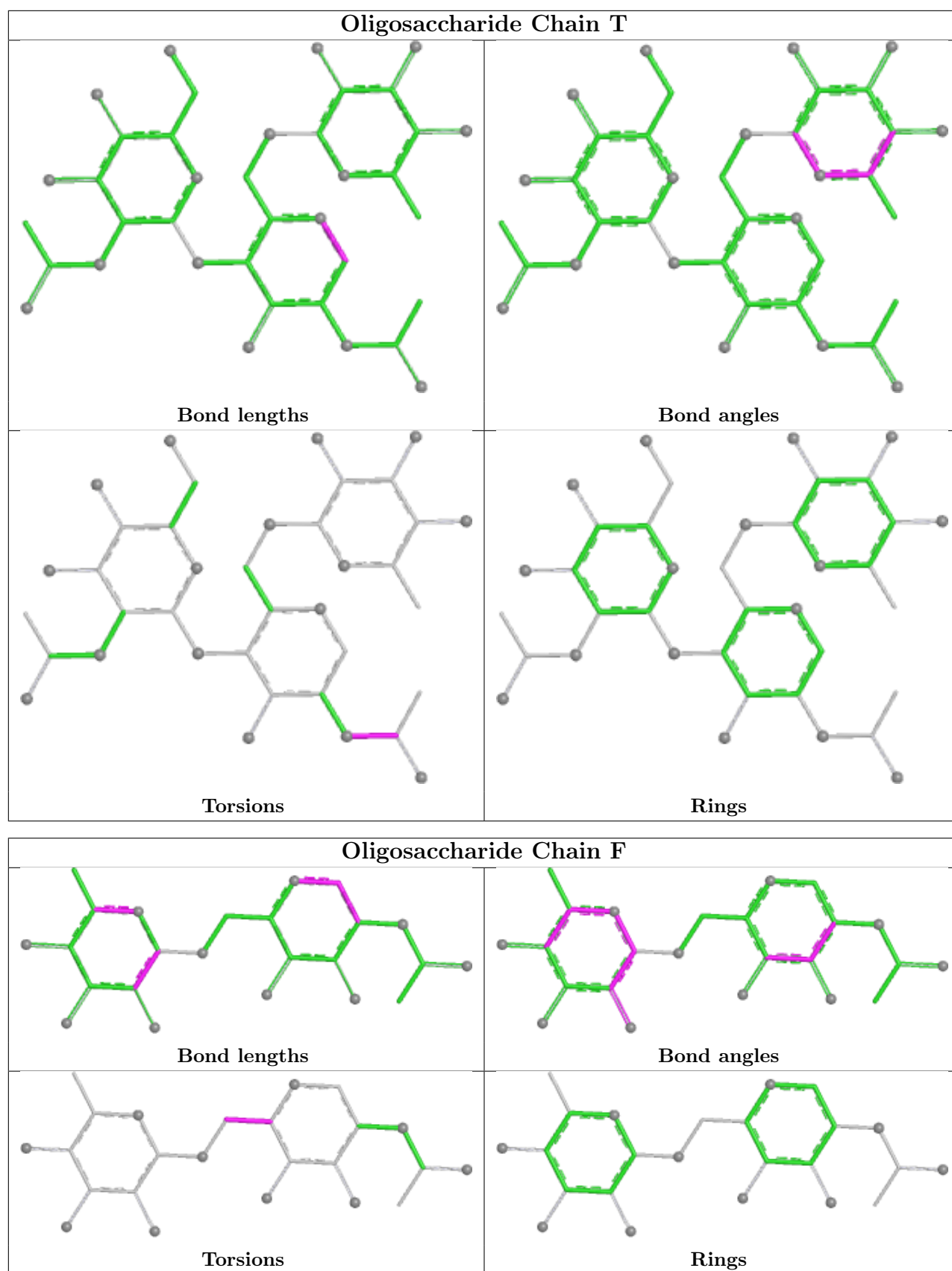
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	M	3	MAN	1	0
2	O	1	NAG	1	0
4	N	1	NAG	2	0
4	W	2	NAG	2	0
5	M	2	NAG	1	0
7	R	5	MAN	1	0
3	U	2	FUC	1	0
2	E	1	NAG	1	0
7	R	3	MAN	1	0
7	R	2	NAG	2	0
4	X	1	NAG	1	0
2	J	1	NAG	1	0
4	H	2	NAG	1	0
3	U	1	NAG	1	0
2	O	3	FUC	1	0

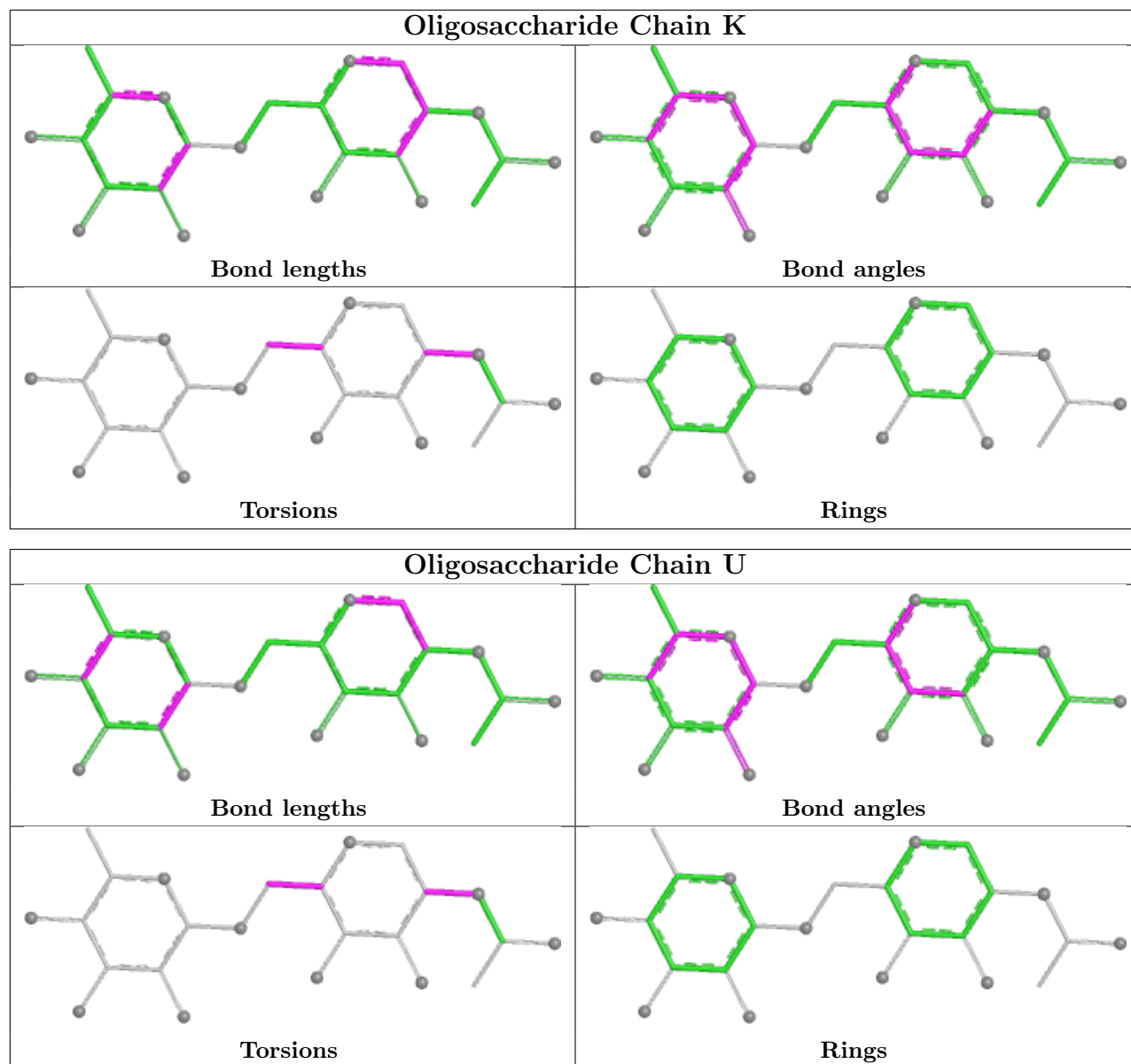
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

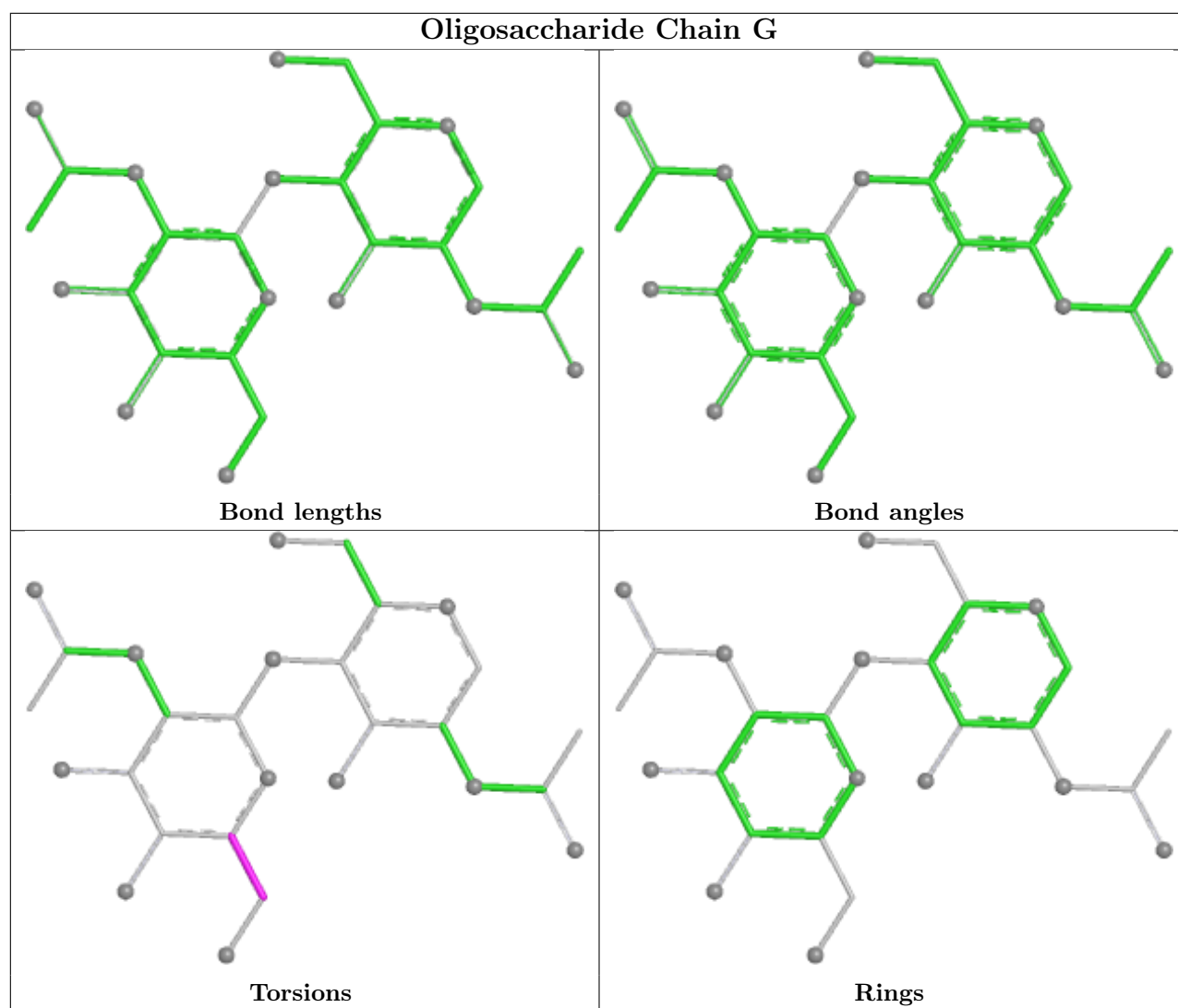


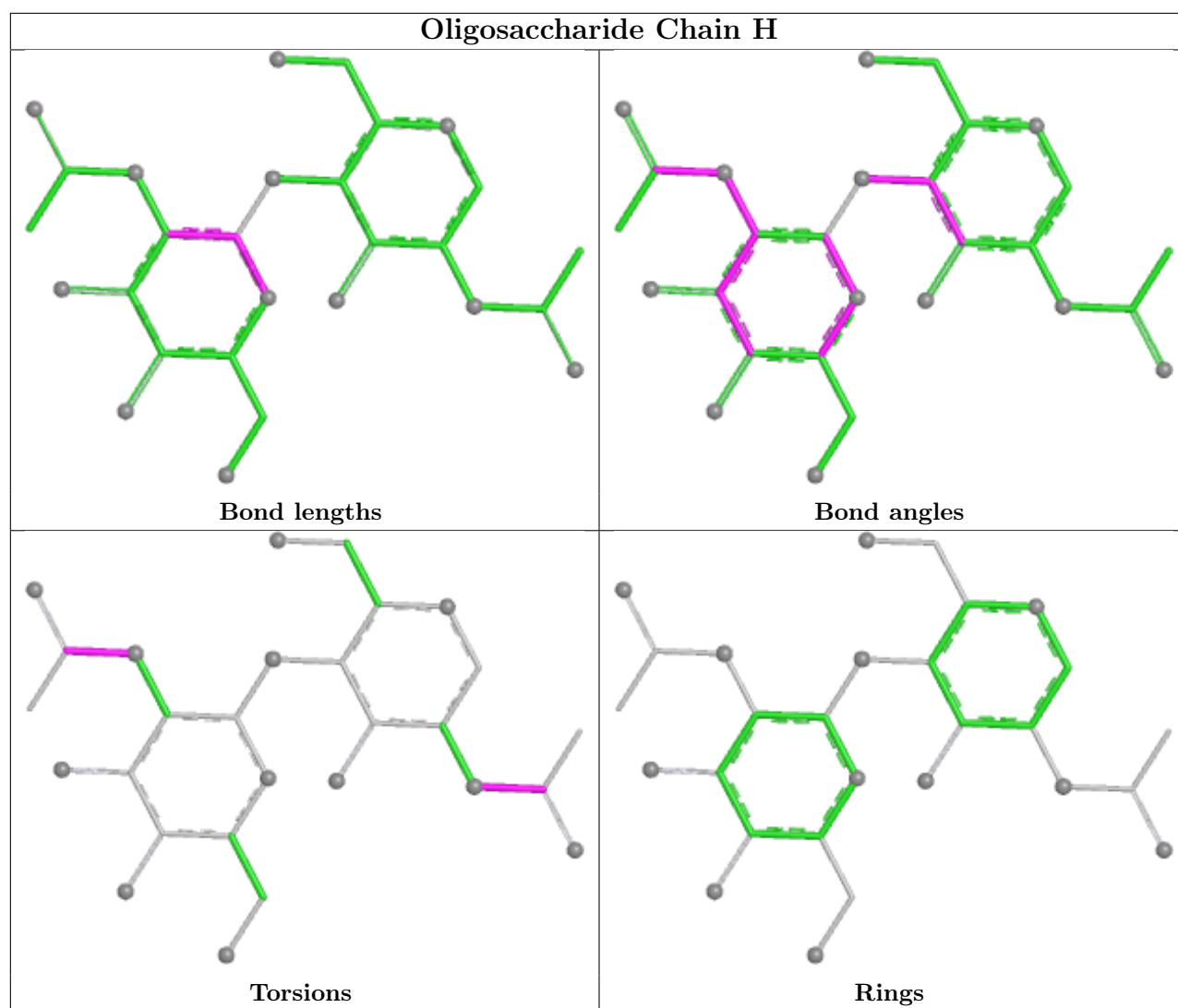


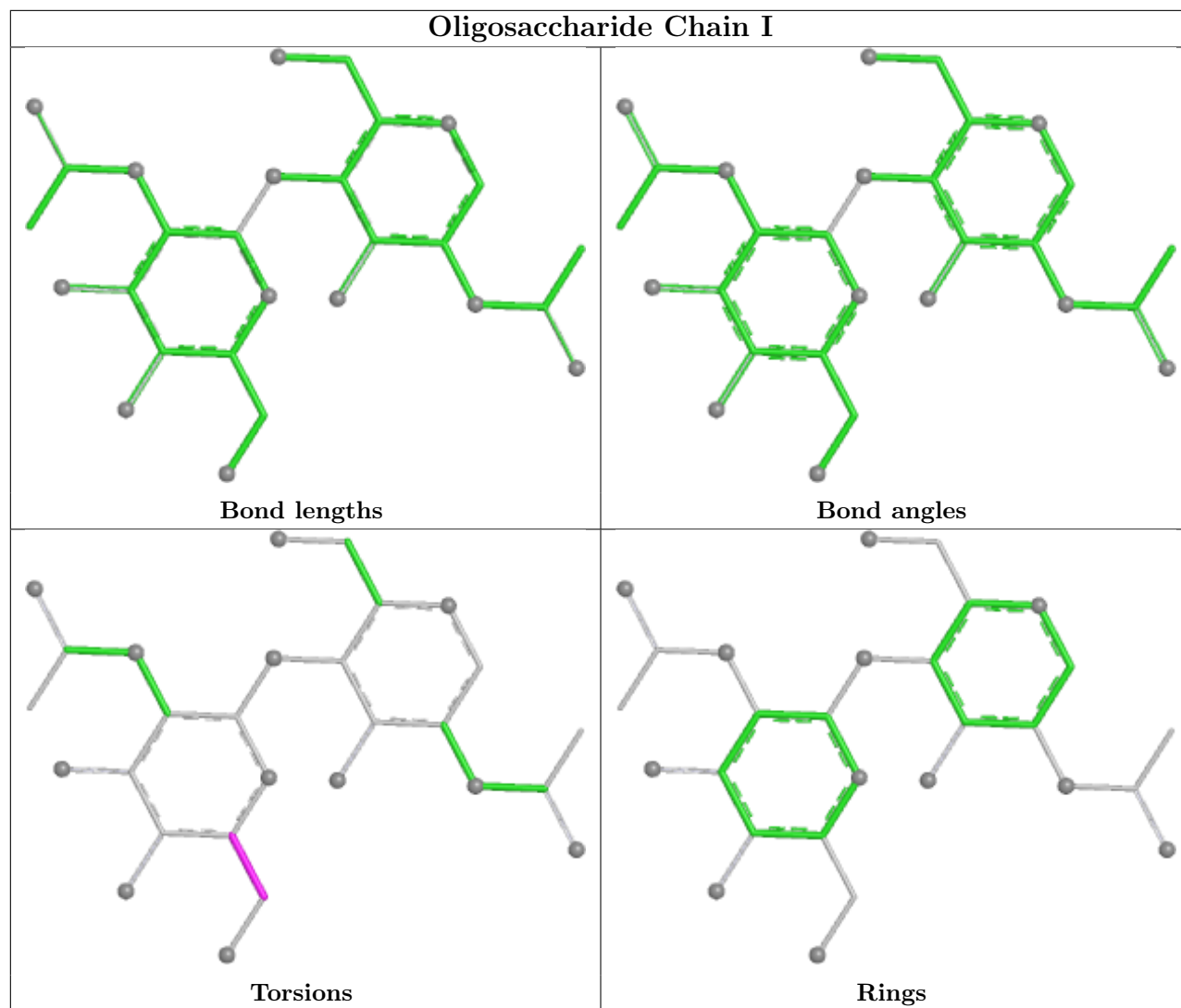


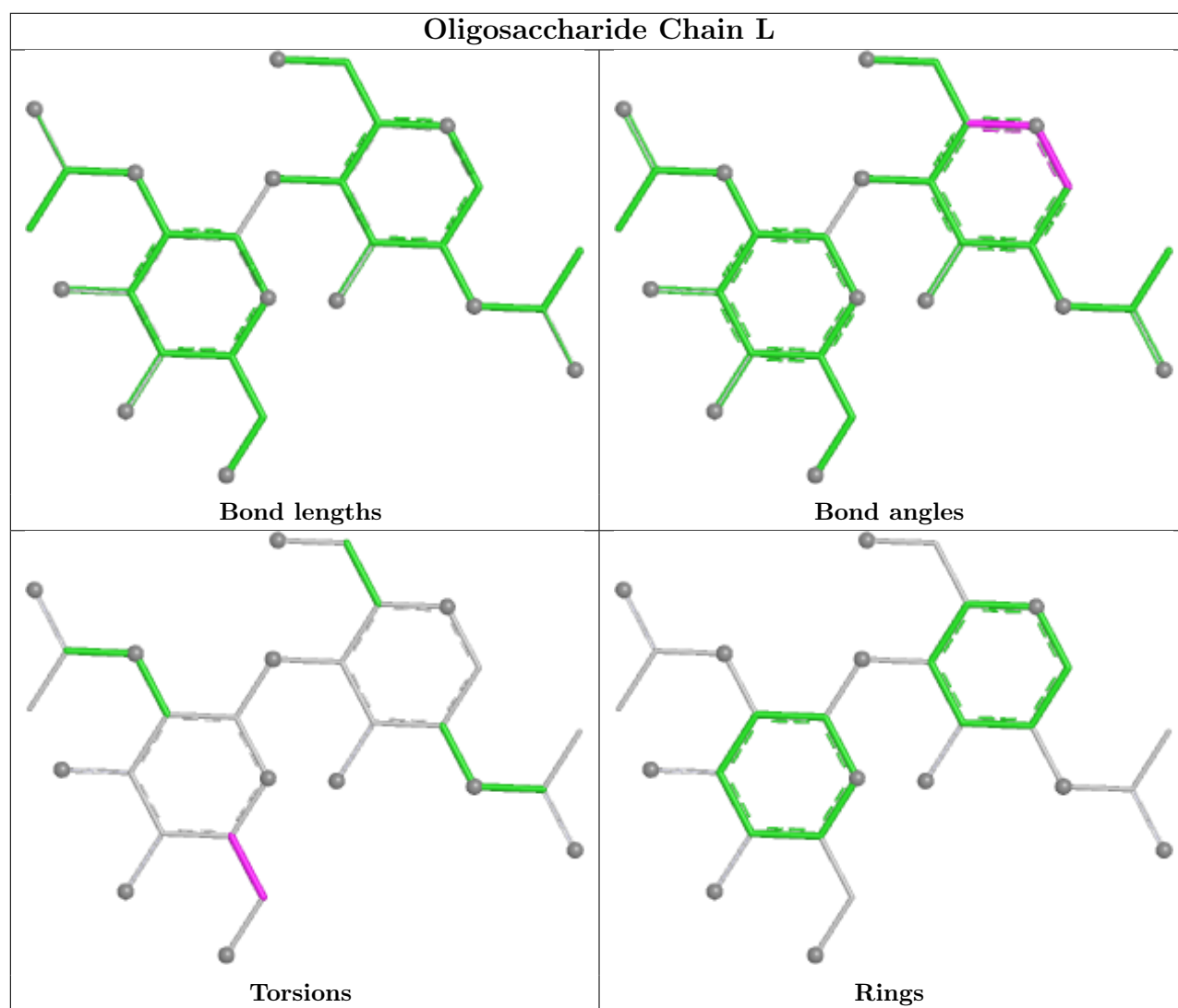


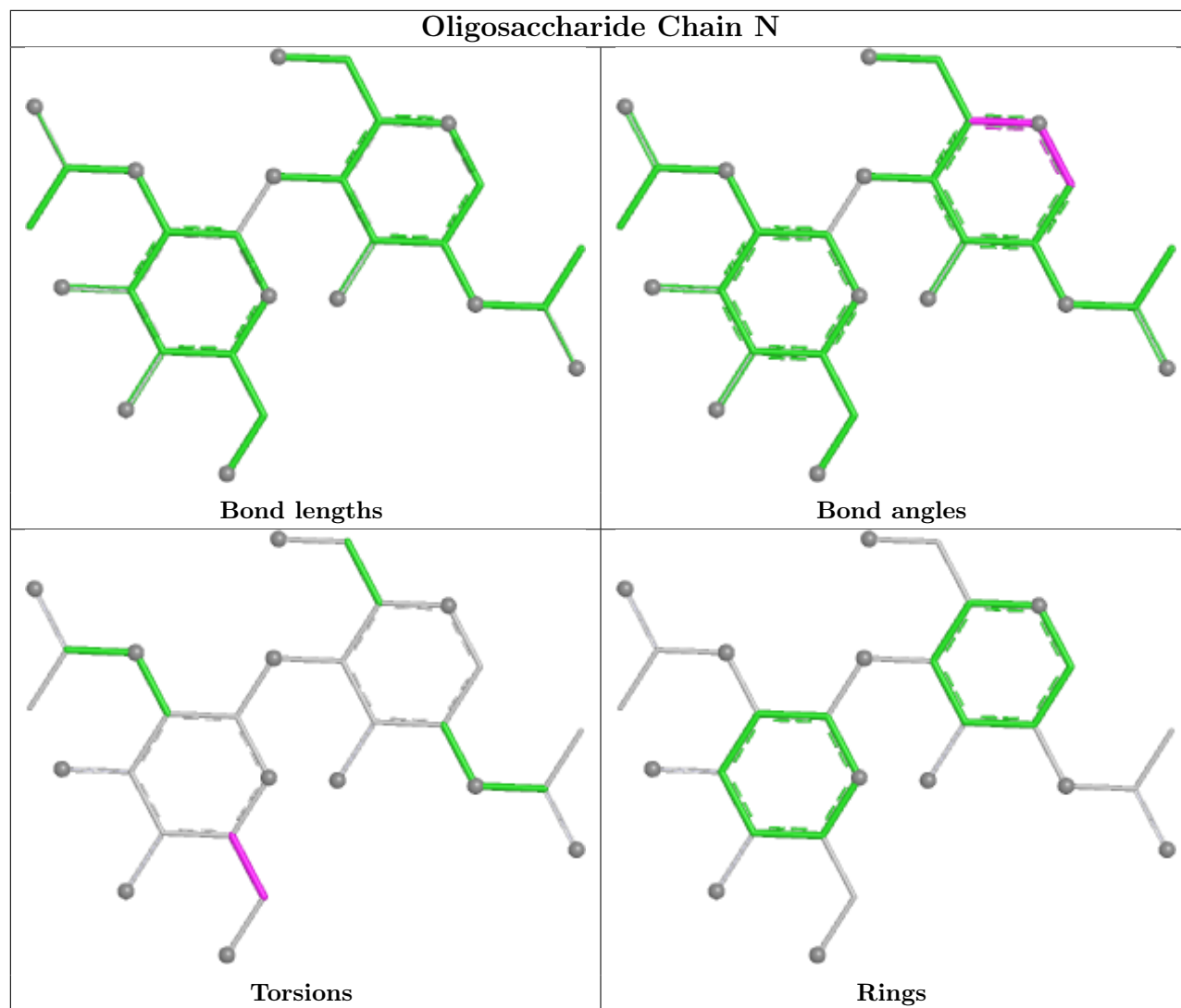


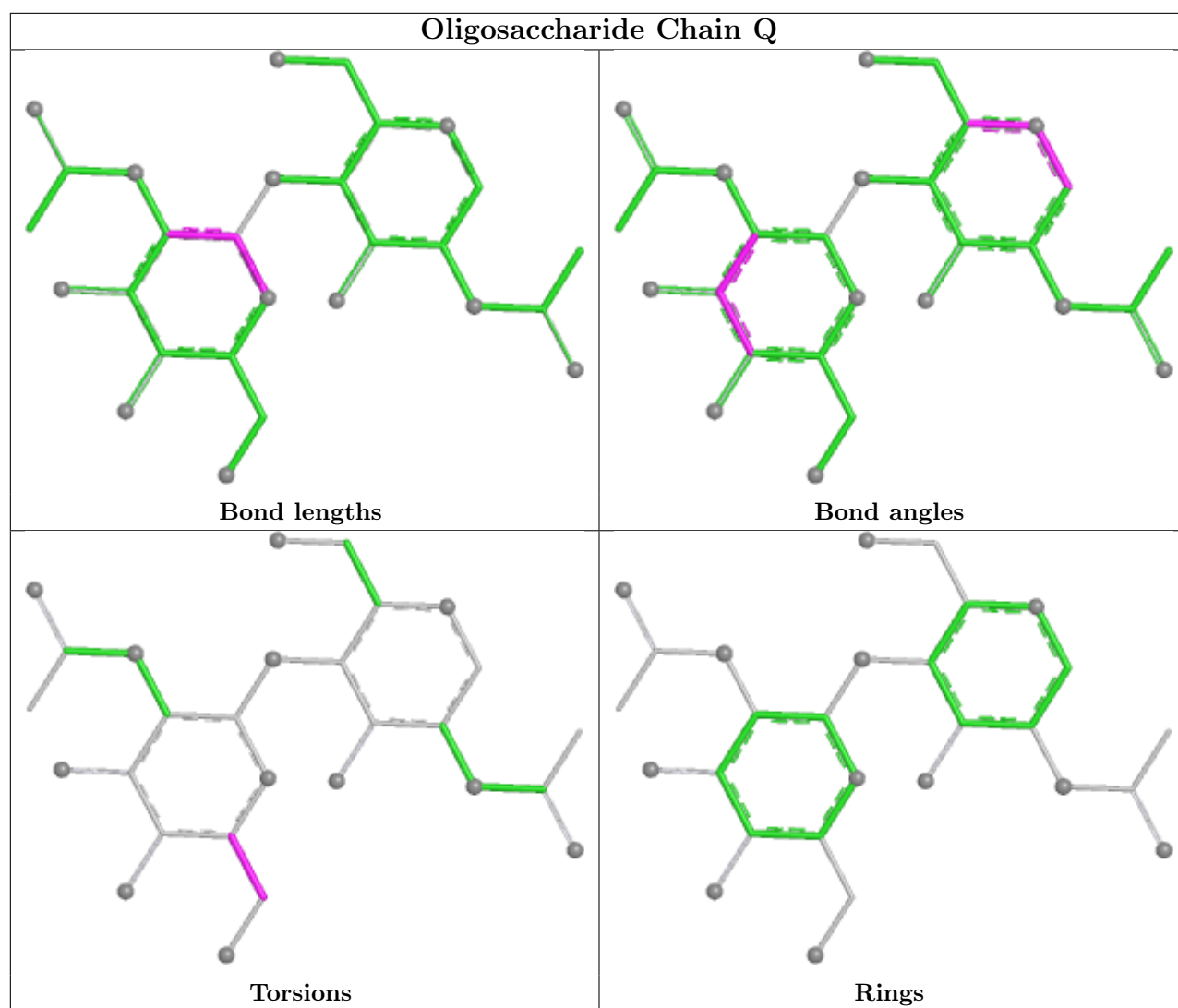


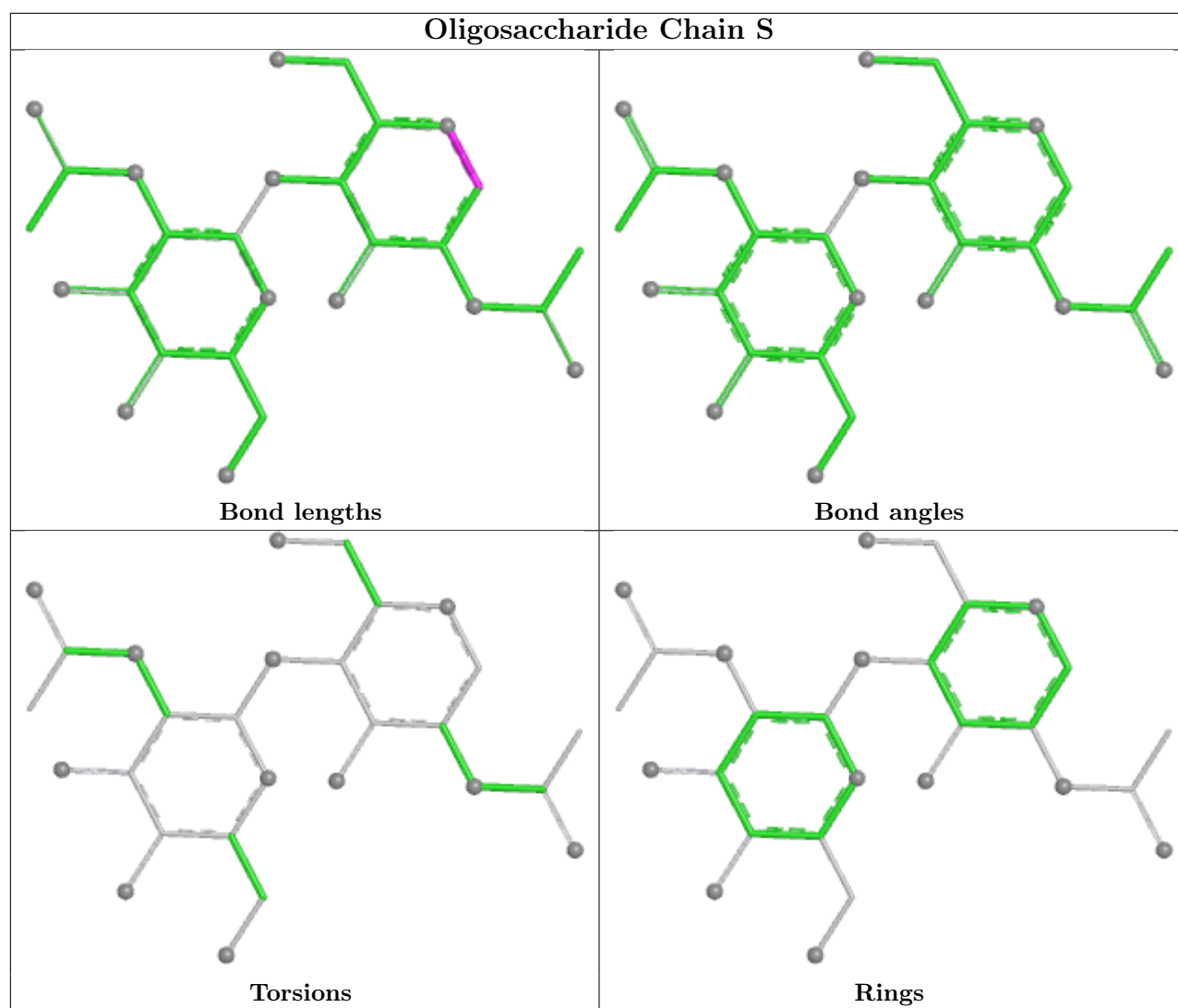


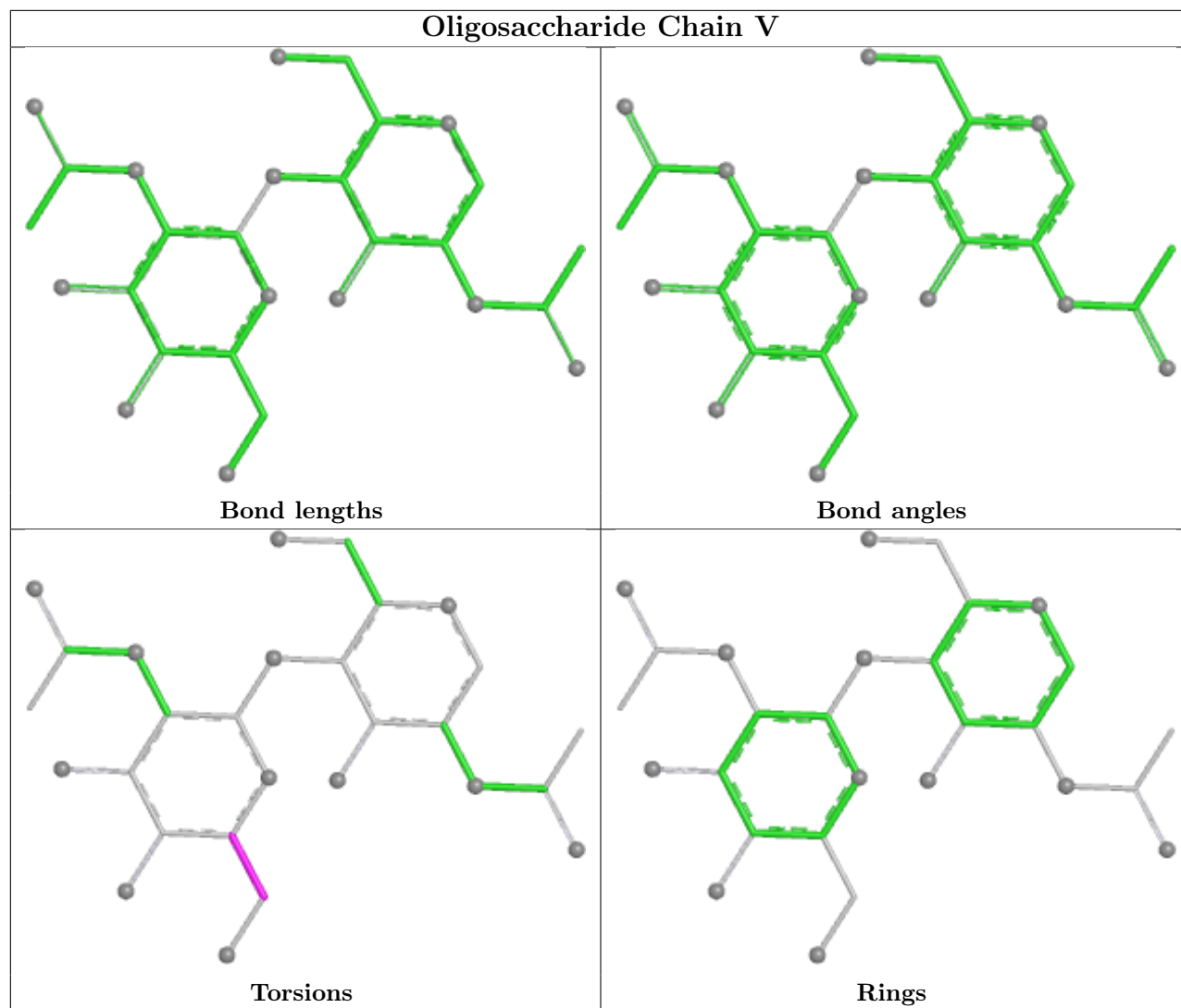


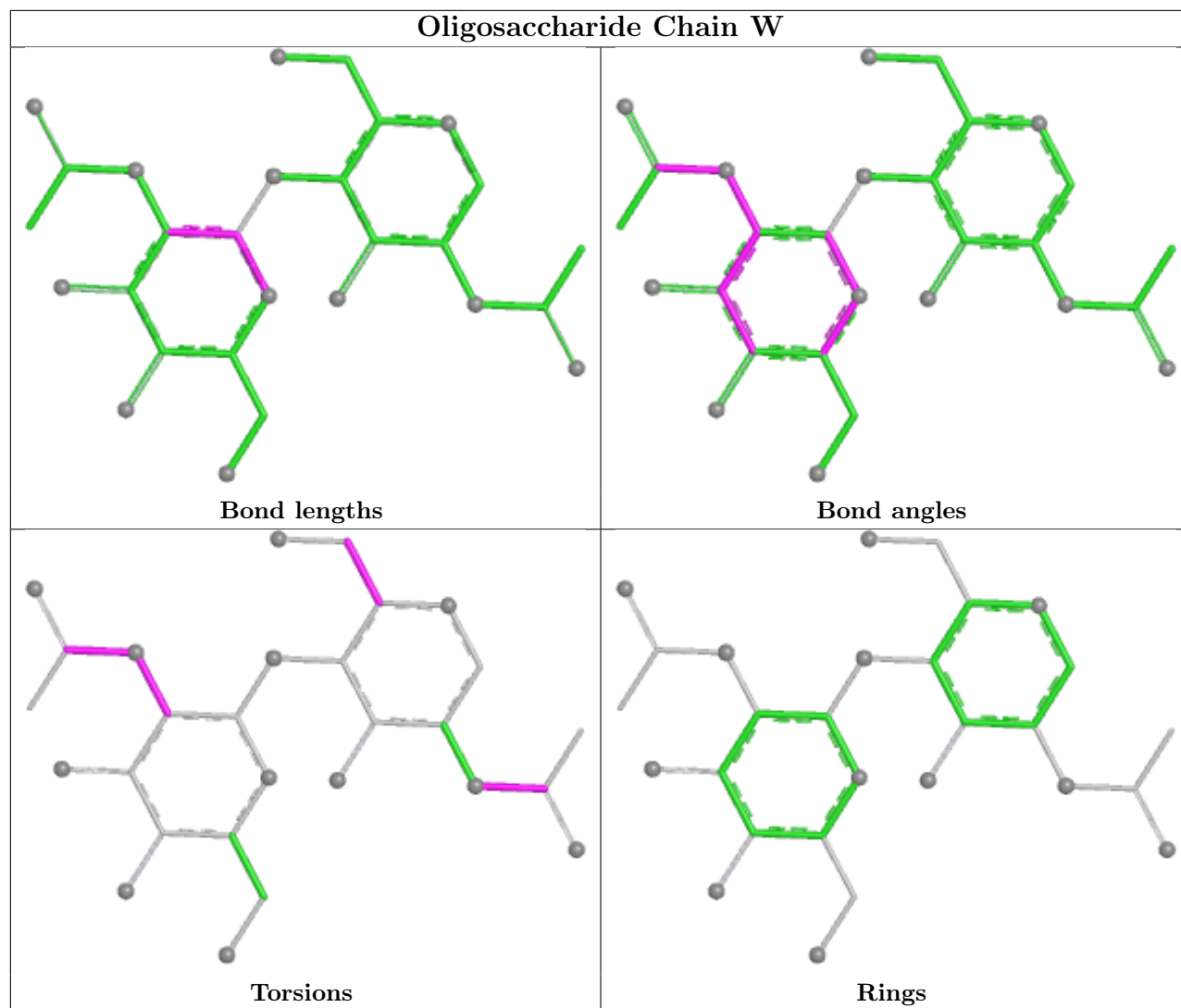


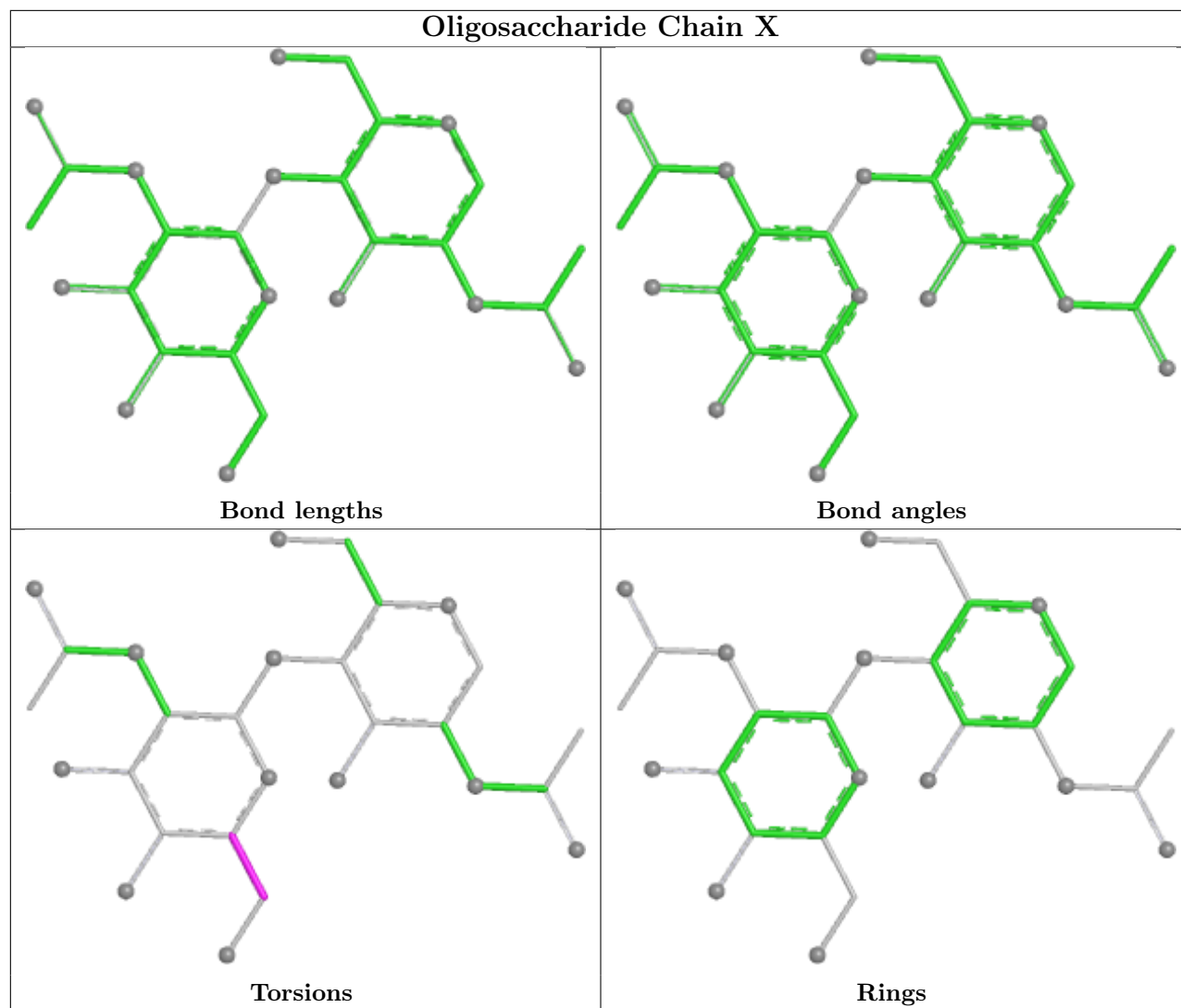




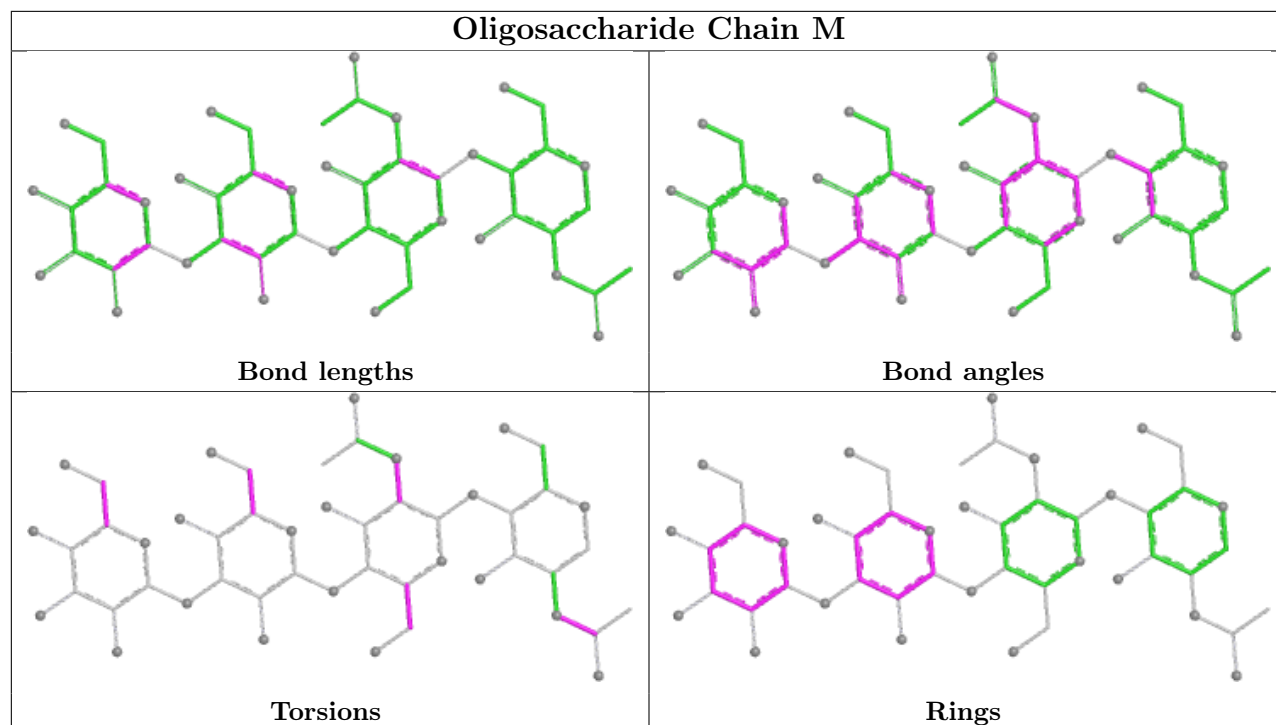




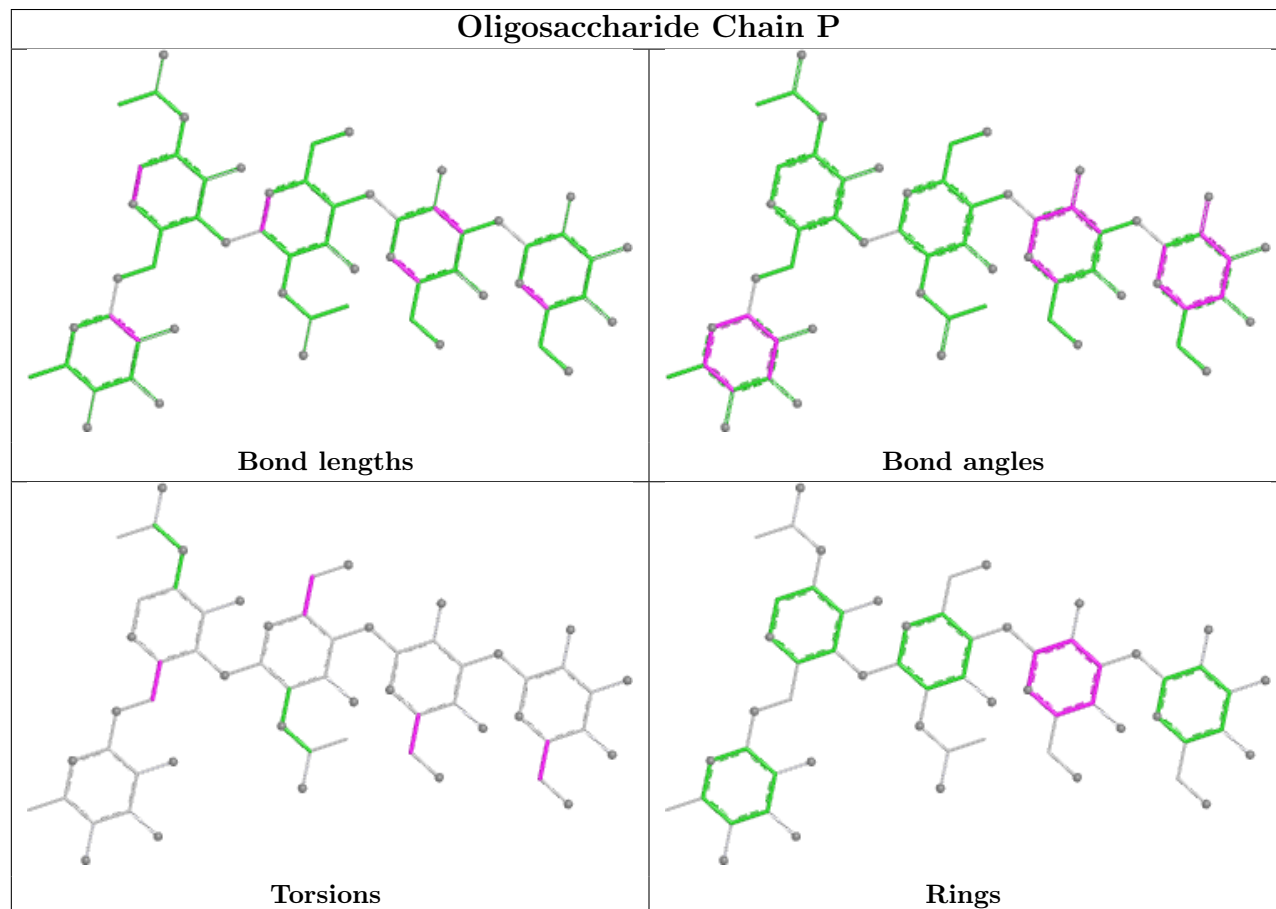


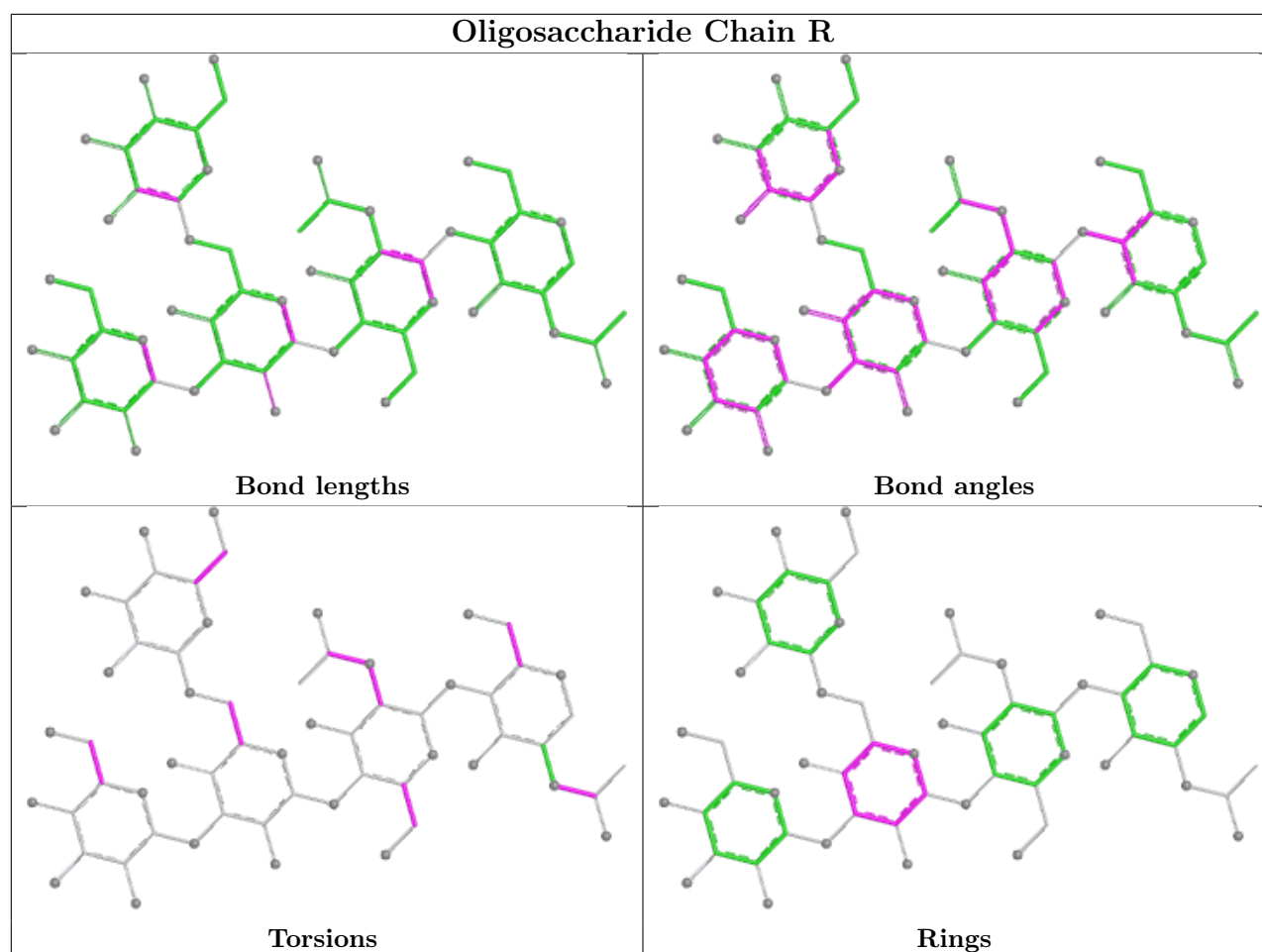


Oligosaccharide Chain M



Oligosaccharide Chain P





5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 9 are monoatomic - leaving 5 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
10	NAG	D	501	1	14,14,15	0.73	1 (7%)	17,19,21	0.72	0
9	KOJ	A	503	-	10,10,10	2.92	5 (50%)	12,13,13	3.55	8 (66%)
9	KOJ	C	503	-	10,10,10	1.90	2 (20%)	12,13,13	3.27	6 (50%)
9	KOJ	D	504	-	10,10,10	1.66	3 (30%)	12,13,13	3.79	8 (66%)
9	KOJ	B	504	-	10,10,10	1.69	2 (20%)	12,13,13	2.99	6 (50%)

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	NAG	D	501	1	-	4/6/23/26	0/1/1/1
9	KOJ	A	503	-	-	1/2/2/2	0/1/1/1
9	KOJ	C	503	-	-	0/2/2/2	0/1/1/1
9	KOJ	D	504	-	-	2/2/2/2	0/1/1/1
9	KOJ	B	504	-	-	0/2/2/2	0/1/1/1

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	503	KOJ	O5-C1	5.29	1.47	1.37
9	A	503	KOJ	O5-C5	5.05	1.46	1.35
9	C	503	KOJ	C4-C3	-4.16	1.33	1.44
9	B	504	KOJ	C4-C3	-3.65	1.35	1.44
9	A	503	KOJ	C1-C2	-3.64	1.31	1.34

The worst 5 of 28 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	503	KOJ	O5-C5-C6	7.05	122.54	111.06
9	D	504	KOJ	O5-C1-C2	-6.57	118.74	122.75
9	B	504	KOJ	O5-C5-C6	6.52	121.69	111.06
9	D	504	KOJ	O5-C5-C6	6.44	121.55	111.06
9	C	503	KOJ	O5-C5-C6	6.30	121.32	111.06

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	D	501	NAG	O5-C5-C6-O6
10	D	501	NAG	C1-C2-N2-C7
9	D	504	KOJ	O5-C5-C6-O6
10	D	501	NAG	C3-C2-N2-C7
10	D	501	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	D	504	KOJ	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	446/446 (100%)	-0.38	3 (0%) 84 82	19, 33, 50, 97	0
1	B	446/446 (100%)	-0.12	8 (1%) 67 63	21, 37, 63, 96	0
1	C	446/446 (100%)	-0.24	3 (0%) 84 82	19, 35, 56, 76	0
1	D	446/446 (100%)	0.08	6 (1%) 75 71	23, 43, 68, 104	0
All	All	1784/1784 (100%)	-0.16	20 (1%) 78 75	19, 36, 62, 104	0

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	48	VAL	5.2
1	D	80	PRO	3.9
1	D	470	SER	3.4
1	B	48	VAL	3.4
1	B	118	ARG	3.3

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(Å ²)	Q<0.9
4	NAG	W	2	14/15	0.46	0.23	94,99,103,105	0
4	NAG	H	2	14/15	0.68	0.18	64,81,91,92	0

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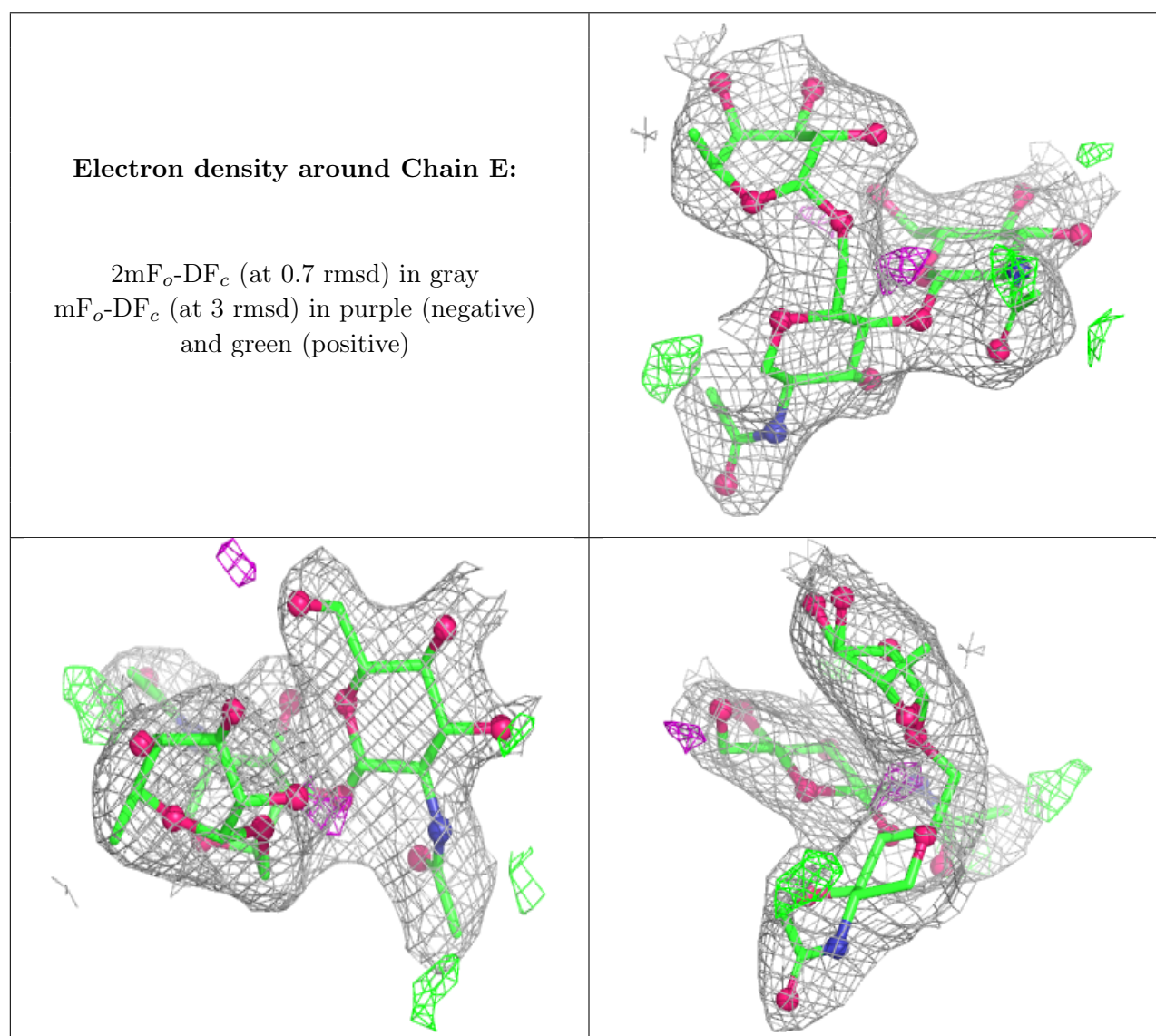
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	K	1	14/15	0.68	0.13	53,66,69,74	0
3	FUC	F	2	10/11	0.71	0.15	62,70,73,73	0
4	NAG	H	1	14/15	0.71	0.16	46,53,74,80	0
4	NAG	X	2	14/15	0.71	0.16	72,75,81,84	0
3	NAG	U	1	14/15	0.72	0.14	49,61,67,69	0
2	NAG	E	2	14/15	0.73	0.14	52,57,64,68	0
4	NAG	V	2	14/15	0.73	0.14	58,72,78,80	0
5	MAN	M	4	11/12	0.73	0.18	55,57,66,70	0
3	NAG	F	1	14/15	0.74	0.14	47,63,69,72	0
2	NAG	T	2	14/15	0.76	0.14	60,70,80,84	0
4	NAG	W	1	14/15	0.77	0.16	68,73,90,92	0
6	MAN	P	4	11/12	0.77	0.14	63,64,70,70	0
2	NAG	J	2	14/15	0.78	0.14	61,73,77,81	0
3	FUC	U	2	10/11	0.78	0.14	54,66,76,81	0
4	NAG	N	2	14/15	0.78	0.14	55,64,71,71	0
2	NAG	T	1	14/15	0.79	0.14	54,59,68,69	0
2	FUC	J	3	10/11	0.79	0.14	53,65,70,72	0
7	NAG	R	2	14/15	0.79	0.20	47,51,58,69	0
7	MAN	R	4	11/12	0.79	0.15	45,53,70,74	0
4	NAG	I	2	14/15	0.80	0.13	50,65,77,83	0
4	NAG	S	2	14/15	0.80	0.14	54,66,73,77	0
2	FUC	T	3	10/11	0.81	0.14	59,72,80,84	0
2	NAG	O	1	14/15	0.81	0.11	49,53,61,63	0
6	MAN	P	3	11/12	0.81	0.12	57,62,66,70	0
5	MAN	M	3	11/12	0.82	0.12	33,41,47,49	0
2	NAG	O	2	14/15	0.82	0.12	52,64,70,74	0
2	NAG	E	1	14/15	0.83	0.11	39,48,54,57	0
2	FUC	O	3	10/11	0.85	0.12	45,53,59,63	0
4	NAG	L	1	14/15	0.85	0.11	34,39,46,49	0
5	NAG	M	2	14/15	0.85	0.12	36,44,51,51	0
3	FUC	K	2	10/11	0.86	0.11	58,71,78,81	0
4	NAG	G	2	14/15	0.86	0.09	43,56,63,68	0
2	NAG	J	1	14/15	0.86	0.11	52,59,66,74	0
7	MAN	R	3	11/12	0.86	0.11	47,49,55,55	0
6	NAG	P	2	14/15	0.86	0.12	53,61,67,74	0
4	NAG	L	2	14/15	0.87	0.10	42,49,58,60	0
4	NAG	Q	2	14/15	0.87	0.12	42,49,57,63	0
4	NAG	X	1	14/15	0.88	0.10	46,58,61,69	0
4	NAG	N	1	14/15	0.88	0.11	35,45,56,63	0
6	FUC	P	5	10/11	0.88	0.11	43,52,55,56	0
7	NAG	R	1	14/15	0.89	0.10	33,39,45,47	0
4	NAG	V	1	14/15	0.89	0.09	47,51,57,69	0

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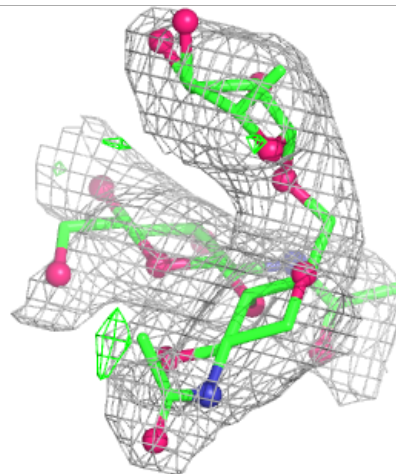
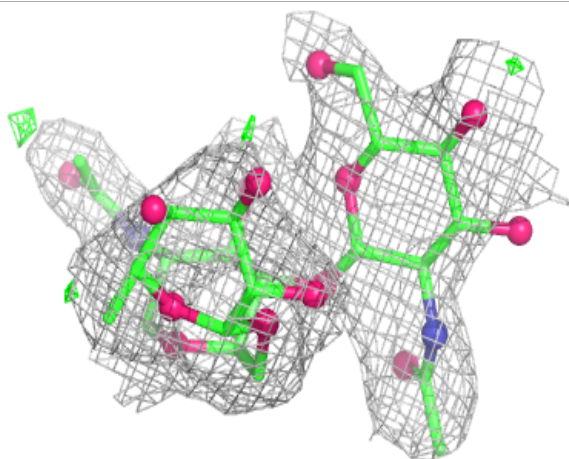
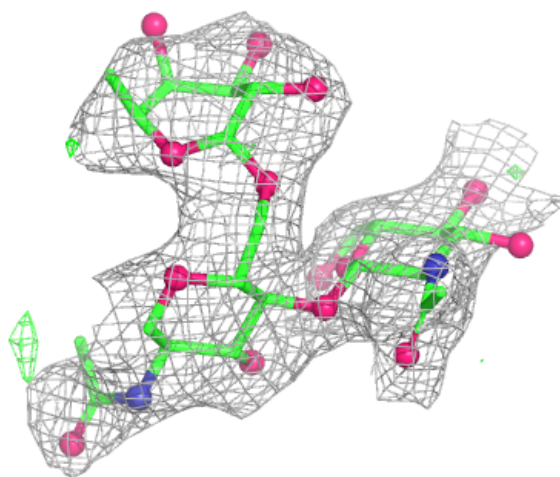
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	MAN	R	5	11/12	0.89	0.11	46,54,58,62	0
6	NAG	P	1	14/15	0.91	0.10	35,44,54,57	0
2	FUC	E	3	10/11	0.91	0.09	38,43,47,56	0
4	NAG	I	1	14/15	0.92	0.08	38,45,50,54	0
4	NAG	S	1	14/15	0.92	0.08	37,47,57,58	0
5	NAG	M	1	14/15	0.93	0.09	31,35,42,42	0
4	NAG	Q	1	14/15	0.93	0.07	30,37,42,42	0
4	NAG	G	1	14/15	0.95	0.07	36,41,47,50	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.



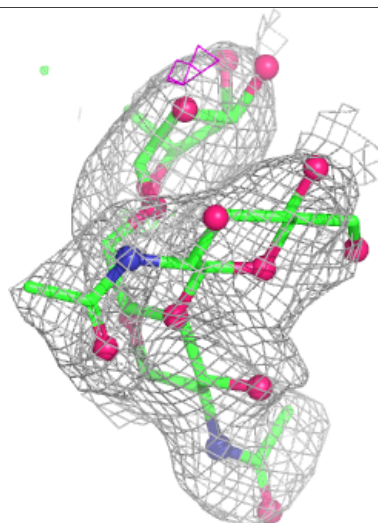
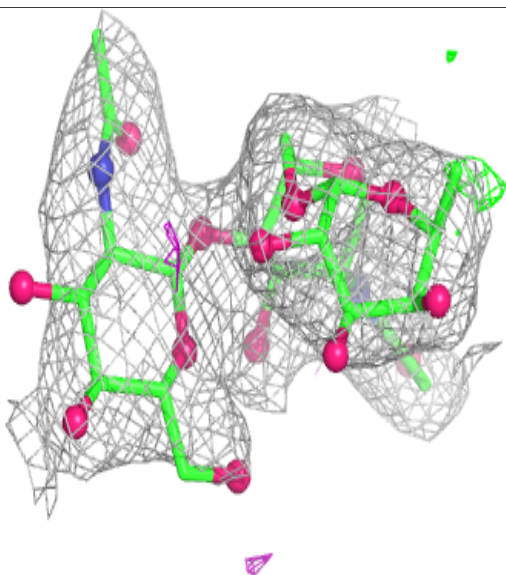
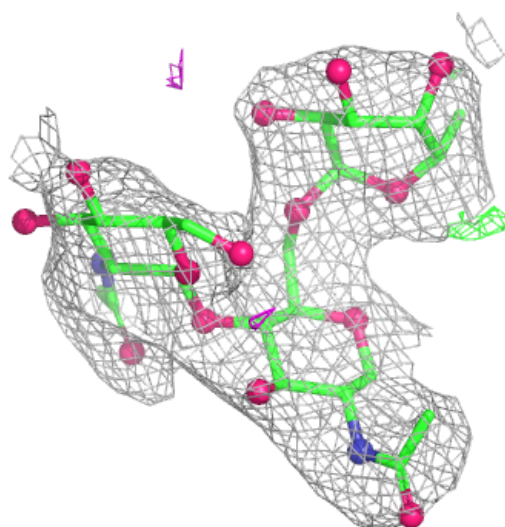
Electron density around Chain J:

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



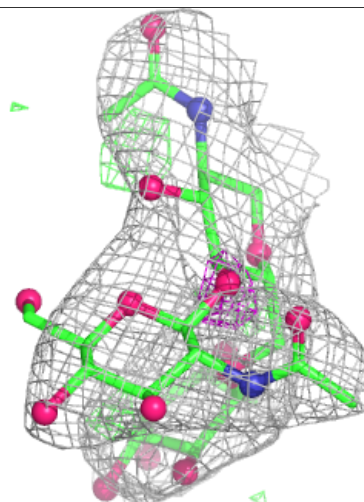
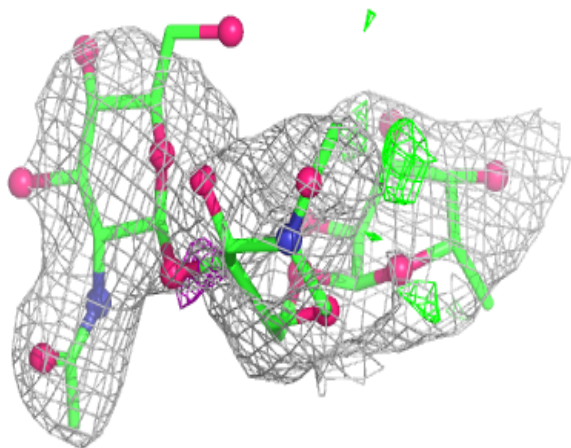
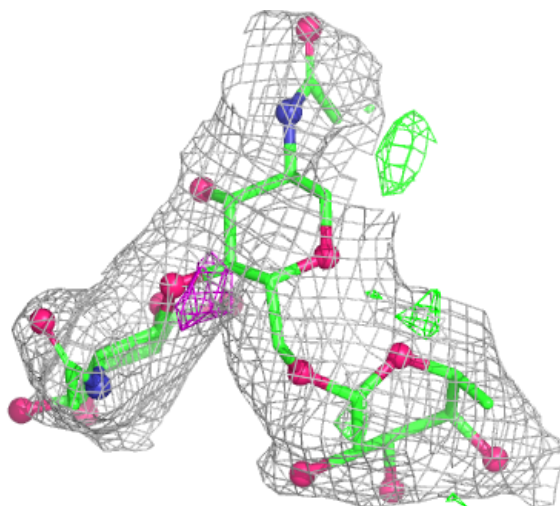
Electron density around Chain O:

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



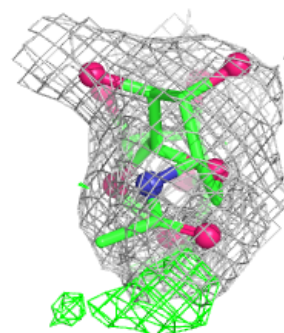
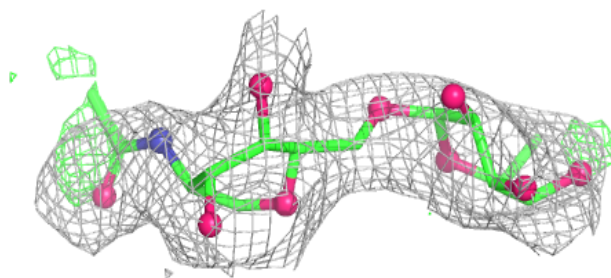
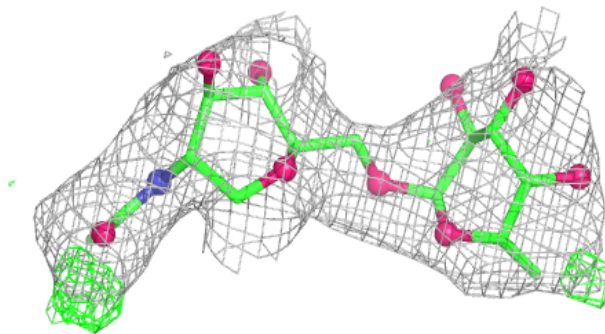
Electron density around Chain T:

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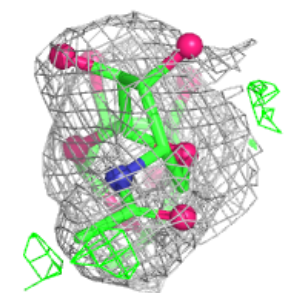
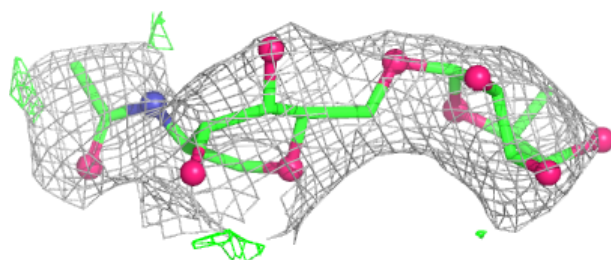
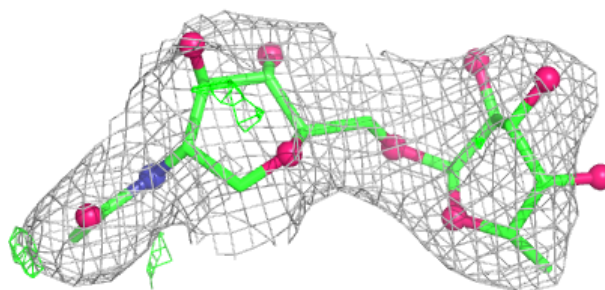


Electron density around Chain F:

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and green (positive)

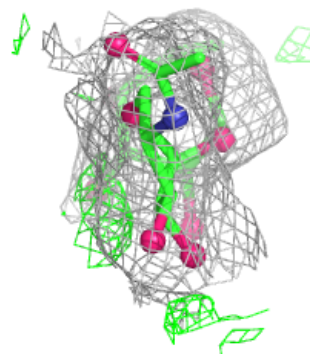
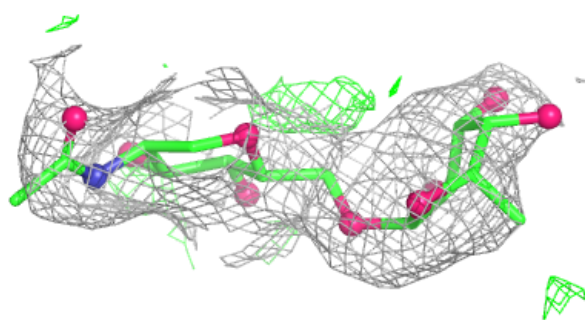
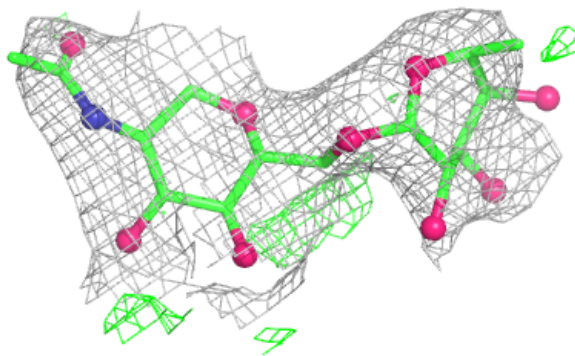
**Electron density around Chain K:**

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



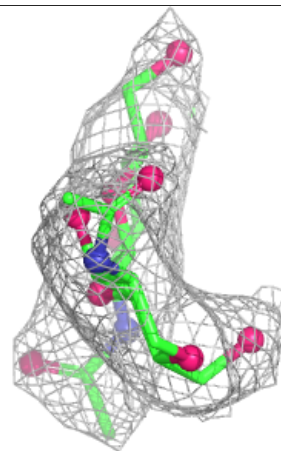
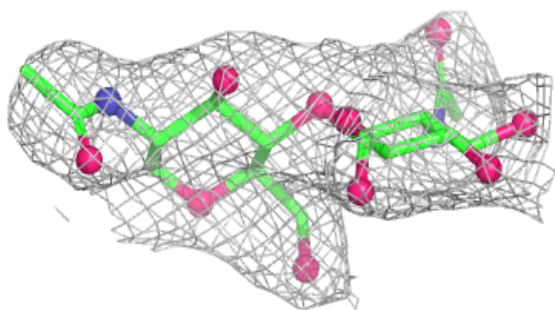
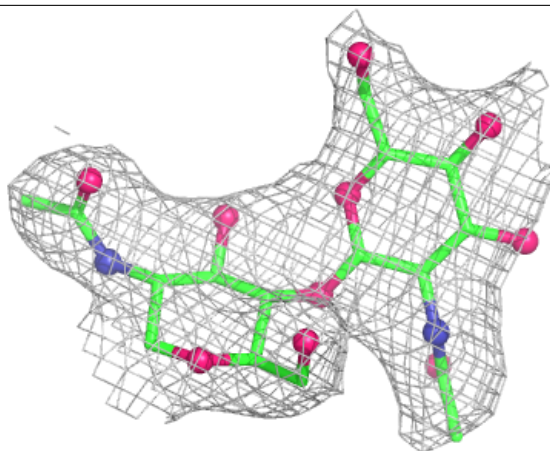
Electron density around Chain U:

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



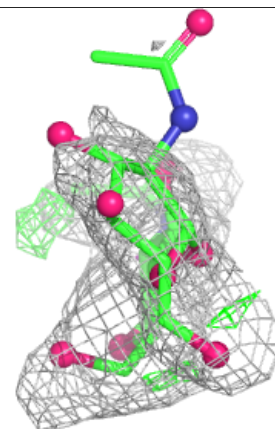
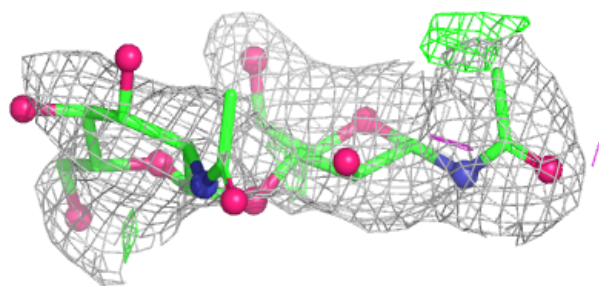
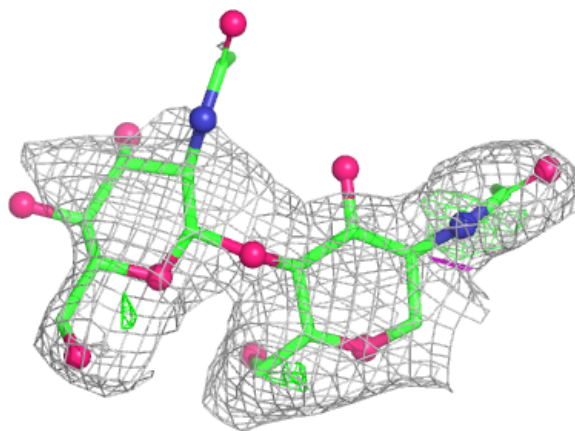
Electron density around Chain G:

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

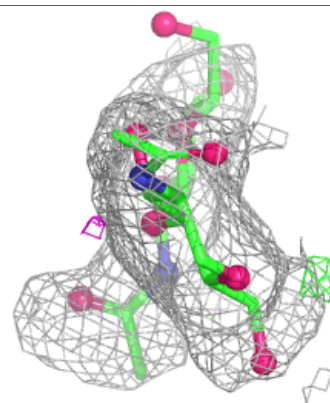
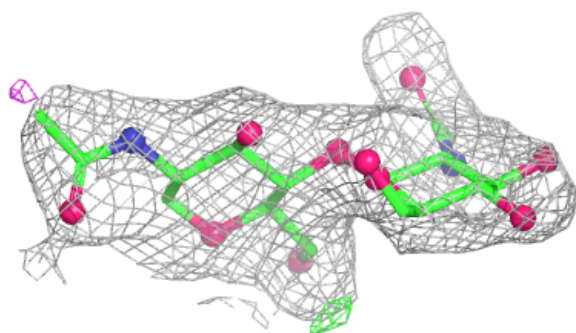
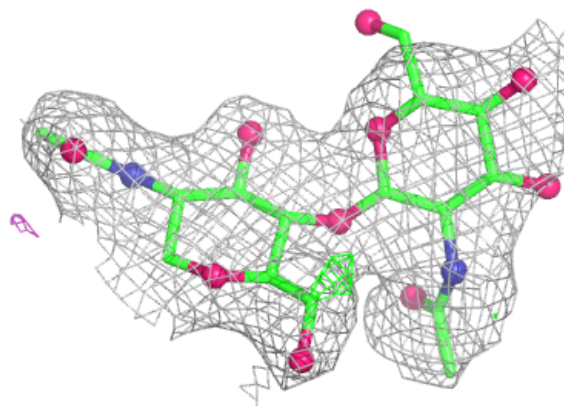


Electron density around Chain H:

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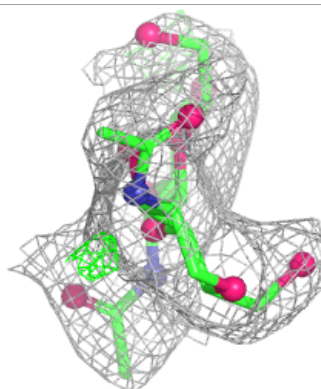
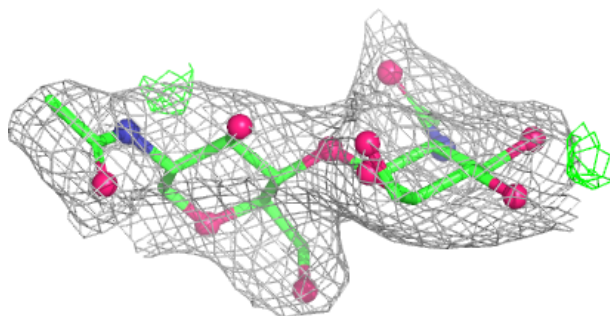
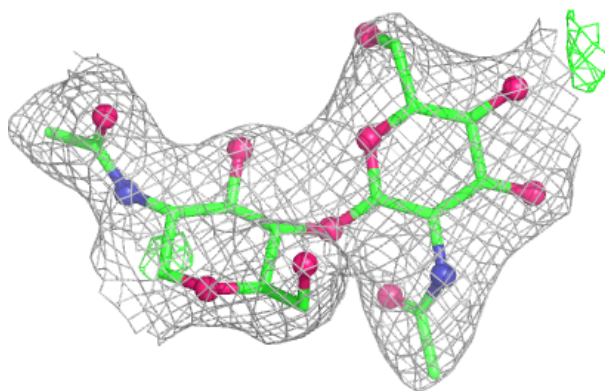
**Electron density around Chain I:**

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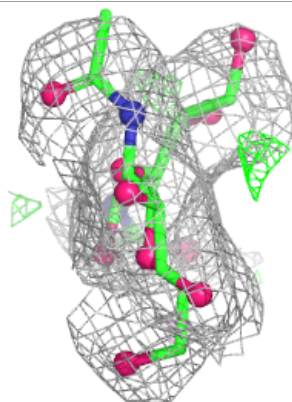
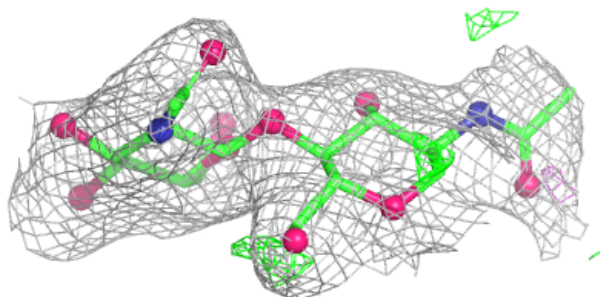
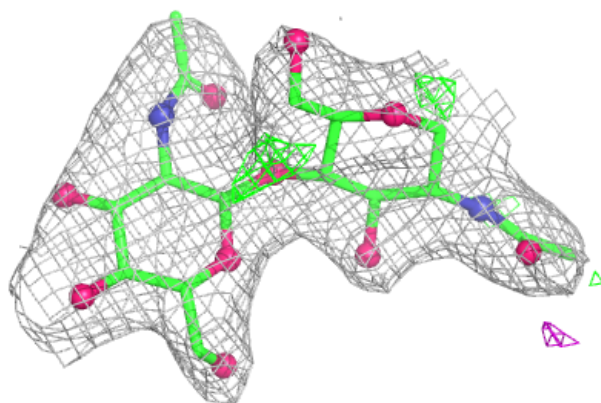


Electron density around Chain L:

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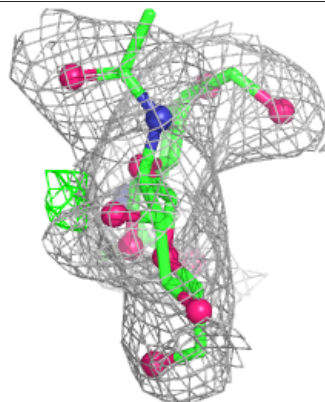
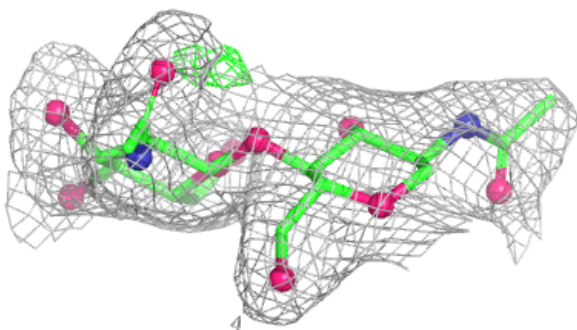
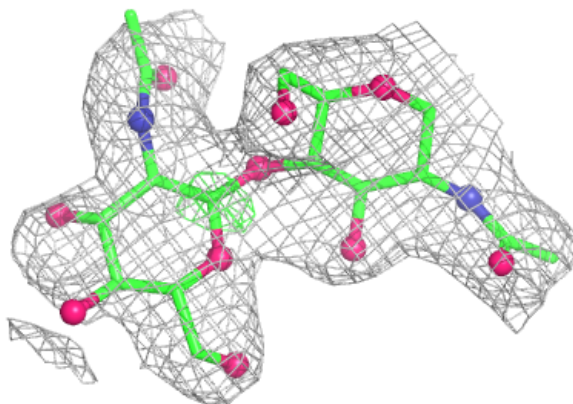
**Electron density around Chain N:**

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and green (positive)



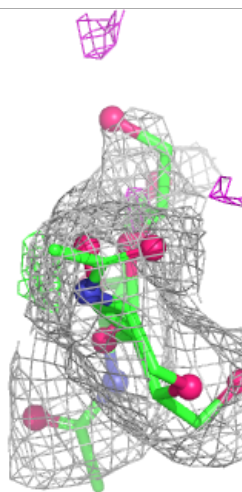
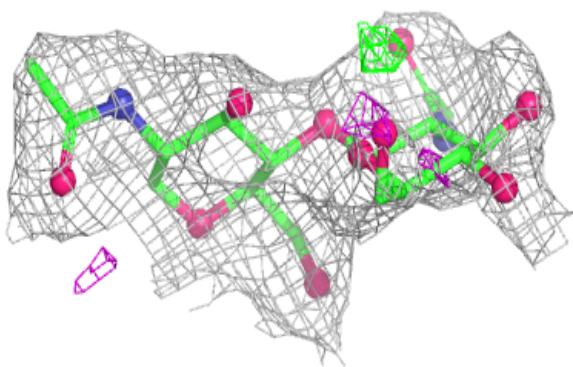
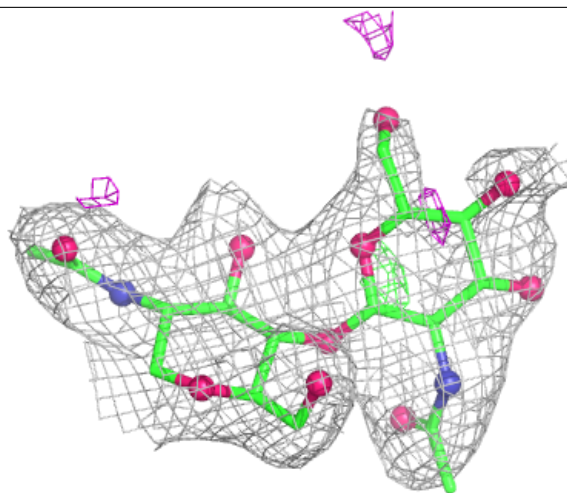
Electron density around Chain Q:

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and green (positive)



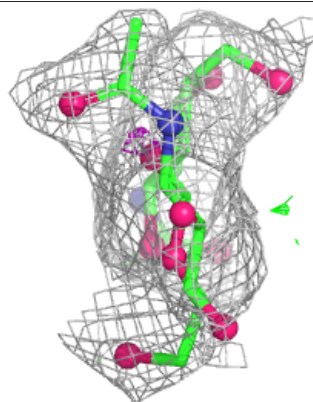
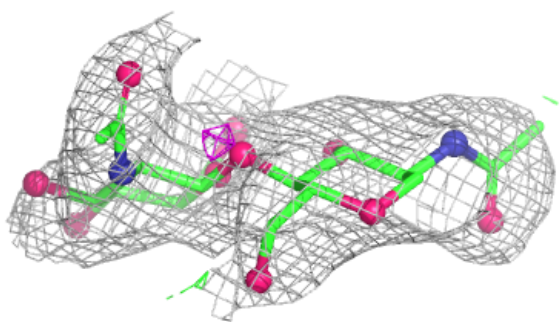
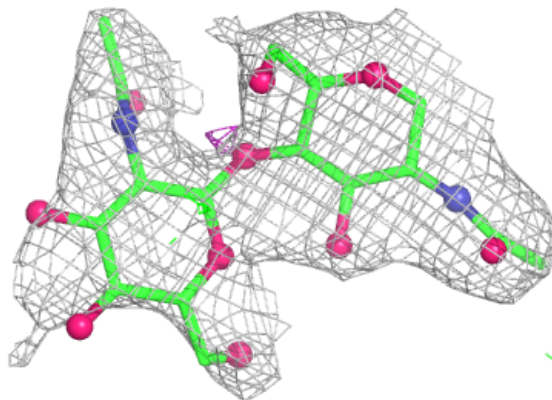
Electron density around Chain S:

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and green (positive)



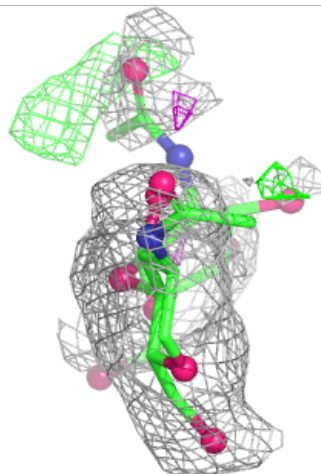
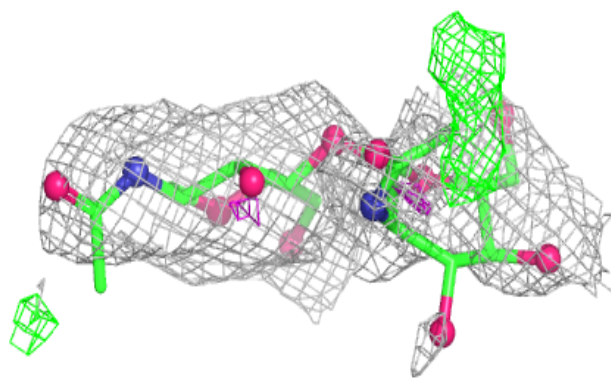
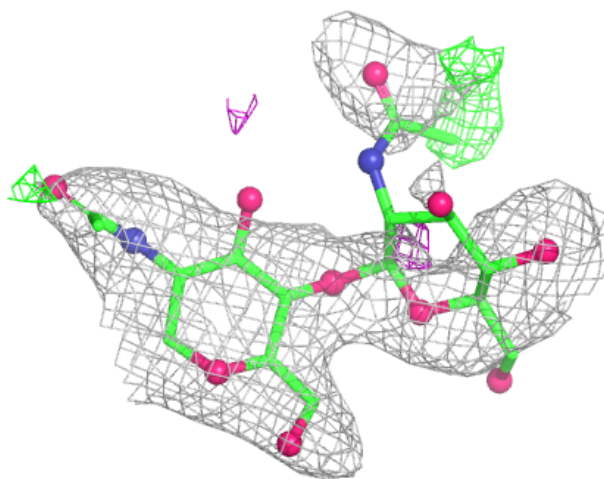
Electron density around Chain V:

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



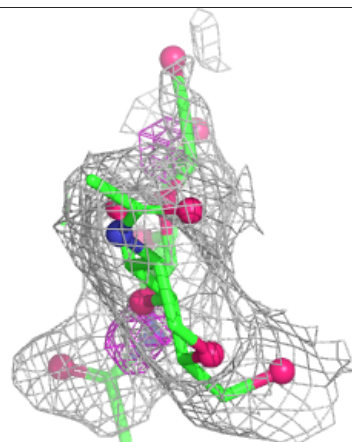
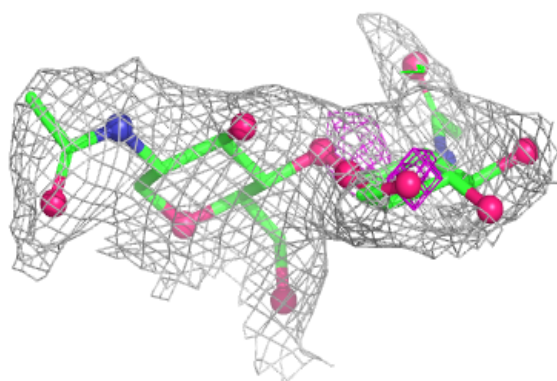
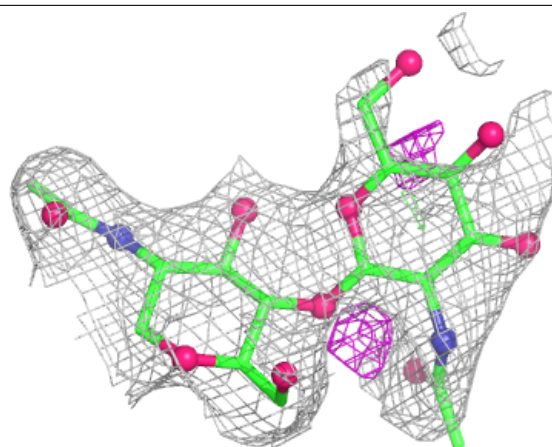
Electron density around Chain W:

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and green (positive)

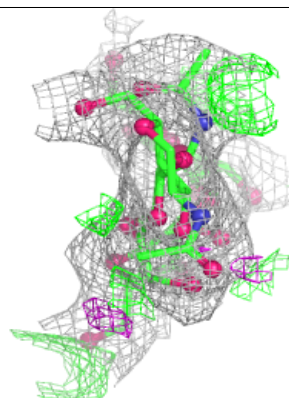
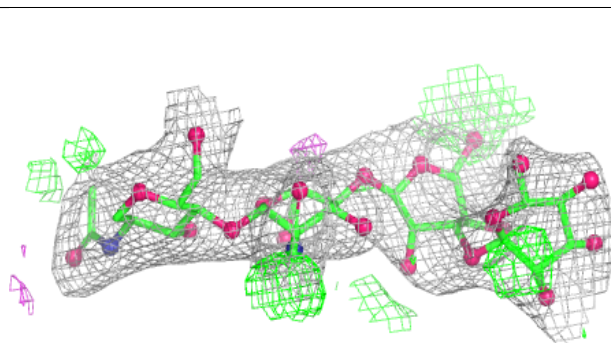
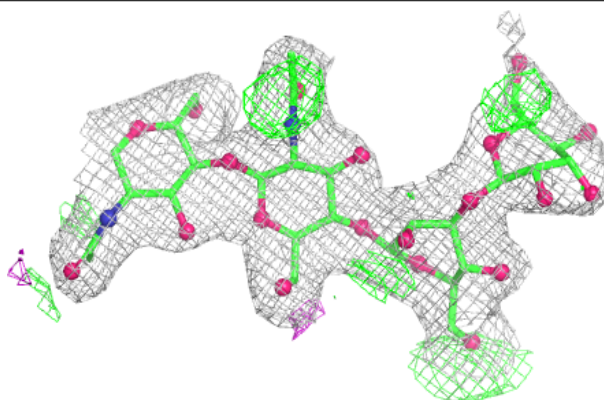


Electron density around Chain X:

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

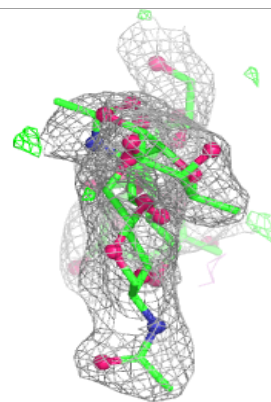
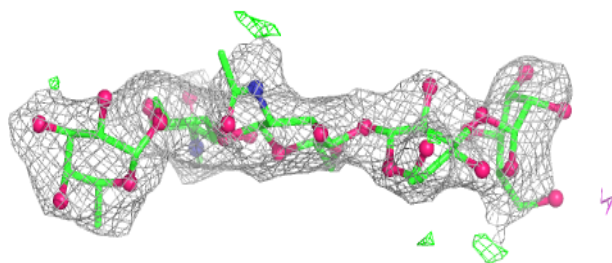
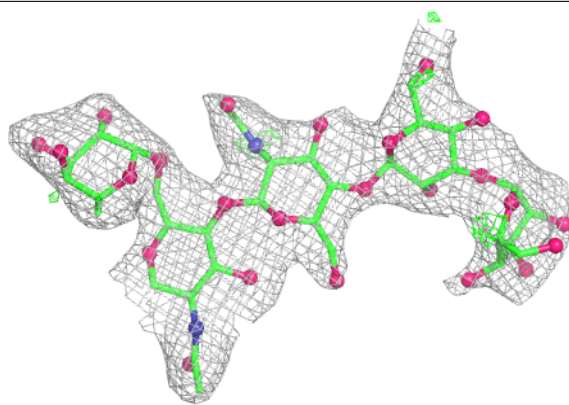
**Electron density around Chain M:**

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

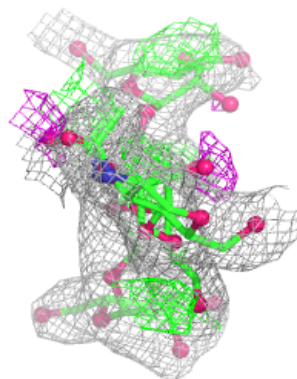
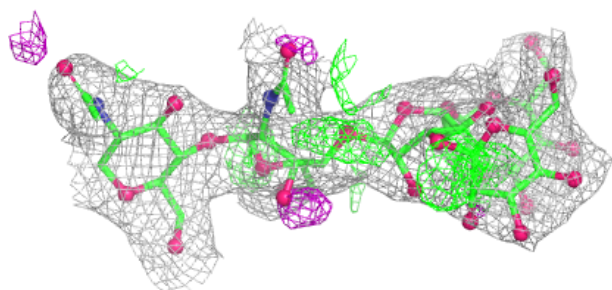
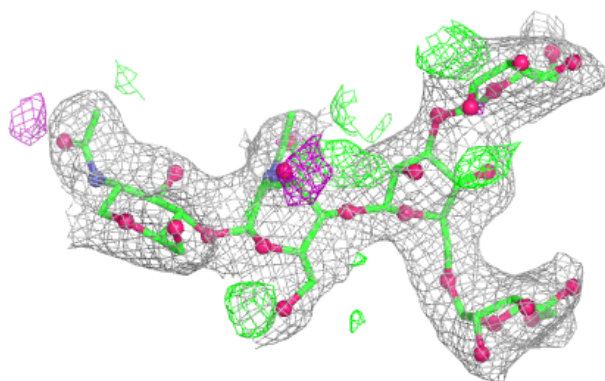


Electron density around Chain P:

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

**Electron density around Chain R:**

$2mF_o-DF_c$ (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	NAG	D	501	14/15	0.71	0.17	50,61,71,76	0
9	KOJ	D	504	10/10	0.85	0.14	52,58,60,64	0
9	KOJ	B	504	10/10	0.88	0.13	29,37,42,45	0
9	KOJ	C	503	10/10	0.89	0.13	35,41,44,44	0
9	KOJ	A	503	10/10	0.93	0.10	32,38,45,46	0
8	ZN	B	501	1/1	0.97	0.10	95,95,95,95	0
8	ZN	D	502	1/1	0.98	0.04	50,50,50,50	0
8	ZN	A	501	1/1	0.98	0.03	40,40,40,40	0
8	ZN	D	503	1/1	0.99	0.03	39,39,39,39	0
8	ZN	B	502	1/1	0.99	0.04	40,40,40,40	0
8	ZN	B	503	1/1	0.99	0.03	27,27,27,27	0
8	ZN	C	501	1/1	0.99	0.05	37,37,37,37	0
8	ZN	C	502	1/1	0.99	0.07	39,39,39,39	0
8	ZN	A	502	1/1	0.99	0.04	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.