



Full wwPDB EM Validation Report ⓘ

Mar 25, 2026 – 02:17 AM UTC

PDB ID : 9JKG / pdb_00009jkg
EMDB ID : EMD-61554
Title : Asymmetric structure of cleaved HIV-1 Tri FPPR envelope glycoprotein trimer
in amphipol-lipid nanodiscs (Tri FPPR.2)
Authors : Qi, Y.; Zhang, S.; Sodroski, J.; Mao, Y.
Deposited on : 2024-09-16
Resolution : 3.50 Å(reported)

This is a Full wwPDB EM Validation 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/EMValidationReportHelp>
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:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

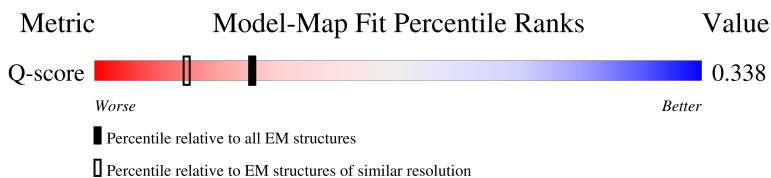
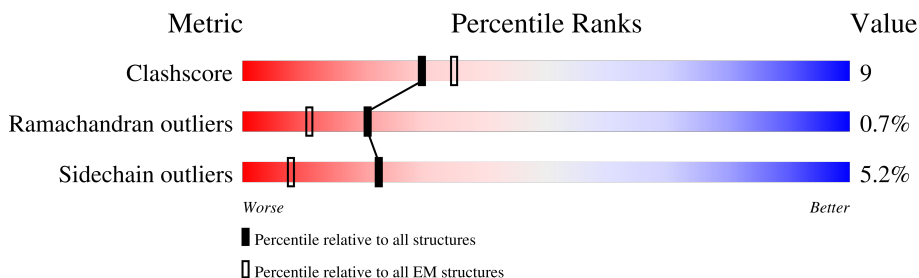
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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.



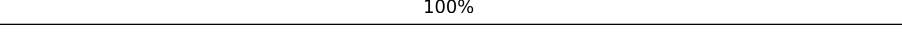
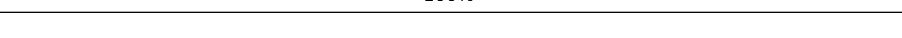
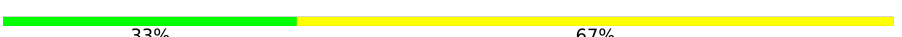
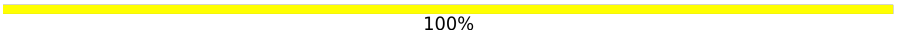
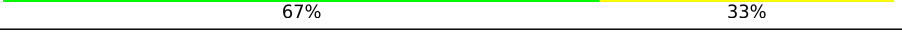
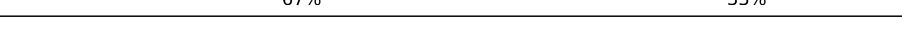
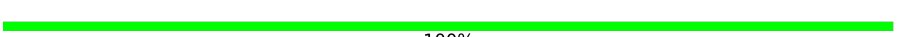
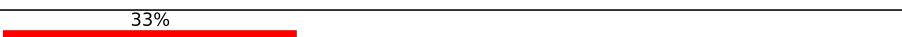
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13950 (3.00 - 4.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	722	
1	B	722	
1	C	722	
1	D	722	

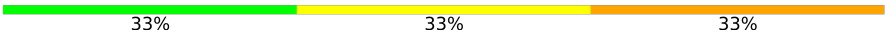
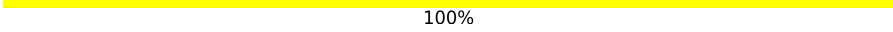
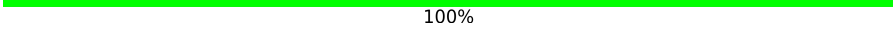
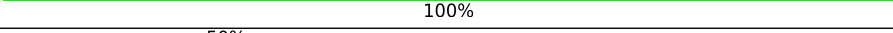
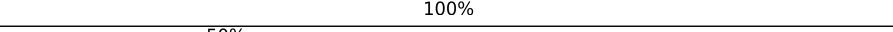
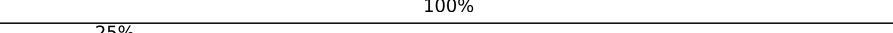
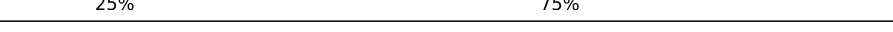
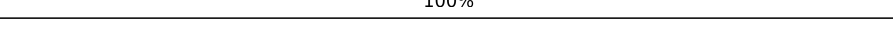
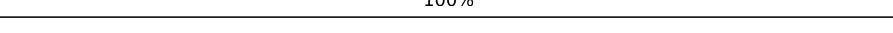
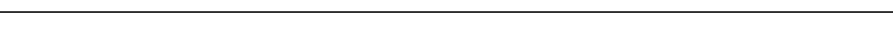
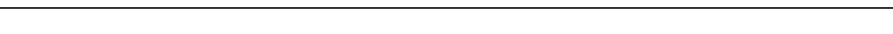
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Mol	Chain	Length	Quality of chain
1	E	722	
1	F	722	
2	0	3	
2	1	3	
2	4	3	
2	8	3	
2	AA	3	
2	BA	3	
2	G	3	
2	I	3	
2	N	3	
2	R	3	
2	T	3	
2	Y	3	
2	Z	3	
2	c	3	
2	d	3	
2	g	3	
2	j	3	
2	k	3	
2	n	3	
2	o	3	
2	s	3	
2	t	3	
2	u	3	

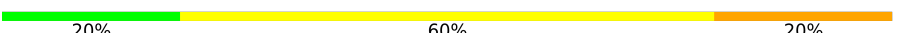
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Mol	Chain	Length	Quality of chain
2	x	3	 67%33%
2	y	3	 33%33%33%
2	z	3	 67%33%
3	EA	2	 50%50%
3	H	2	 50%50%
3	J	2	 100%
3	M	2	 50%100%
3	O	2	 50%50%
3	P	2	 50%50%
3	S	2	 100%
3	b	2	 50%100%
3	h	2	 50%50%
3	v	2	 100%
4	5	4	 25%75%
4	K	4	 75%25%
4	i	4	 100%
4	p	4	 100%
5	3	4	 75%25%
5	9	4	 50%25%25%
5	L	4	 50%50%
5	U	4	 50%50%
5	X	4	 25%50%
5	m	4	 75%25%
5	r	4	 100%
6	2	6	 50%17%33%

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Mol	Chain	Length	Quality of chain
6	Q	6	 67%17%17%
7	6	5	 60%40%
7	V	5	 20%80%
7	e	5	 20%60%20%
7	f	5	 20%80%20%
7	w	5	 20%60%40%
8	W	4	 75%25%
9	CA	4	 75%25%
9	a	4	 25%75%25%
10	l	8	 12%75%25%
11	7	7	 14%43%57%
11	q	7	 29%71%
12	DA	3	 67%100%
12	FA	3	 67%33%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 17991 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Envelope glycoprotein gp160.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	170	Total	C	N	O	S	0	0
			1399	904	235	252	8		
1	C	475	Total	C	N	O	S	0	0
			3751	2344	665	716	26		
1	D	144	Total	C	N	O	S	0	0
			1159	735	200	217	7		
1	E	472	Total	C	N	O	S	0	0
			3722	2328	658	710	26		
1	A	478	Total	C	N	O	S	0	0
			3780	2361	672	721	26		
1	F	144	Total	C	N	O	S	0	0
			1159	735	200	217	7		

There are 216 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	2	MET	-	initiating methionine	UNP G1JZH9
B	3	ARG	-	expression tag	UNP G1JZH9
B	4	VAL	-	expression tag	UNP G1JZH9
B	5	LYS	-	expression tag	UNP G1JZH9
B	6	GLU	-	expression tag	UNP G1JZH9
B	7	LYS	-	expression tag	UNP G1JZH9
B	8	TYR	-	expression tag	UNP G1JZH9
B	9	GLN	-	expression tag	UNP G1JZH9
B	10	HIS	-	expression tag	UNP G1JZH9
B	11	LEU	-	expression tag	UNP G1JZH9
B	12	TRP	-	expression tag	UNP G1JZH9
B	13	ARG	-	expression tag	UNP G1JZH9
B	14	TRP	-	expression tag	UNP G1JZH9
B	15	GLY	-	expression tag	UNP G1JZH9
B	16	TRP	-	expression tag	UNP G1JZH9
B	17	ARG	-	expression tag	UNP G1JZH9
B	18	TRP	-	expression tag	UNP G1JZH9
B	19	GLY	-	expression tag	UNP G1JZH9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	20	THR	-	expression tag	UNP G1JZH9
B	32	THR	VAL	conflict	UNP G1JZH9
B	34	LYS	ASN	conflict	UNP G1JZH9
B	115	GLU	GLN	conflict	UNP G1JZH9
B	532	VAL	ALA	conflict	UNP G1JZH9
B	535	MET	ILE	conflict	UNP G1JZH9
B	543	GLN	LEU	conflict	UNP G1JZH9
B	567	LYS	GLN	conflict	UNP G1JZH9
B	582	THR	ALA	conflict	UNP G1JZH9
B	715	GLY	-	expression tag	UNP G1JZH9
B	716	GLY	-	expression tag	UNP G1JZH9
B	717	GLY	-	expression tag	UNP G1JZH9
B	718	HIS	-	expression tag	UNP G1JZH9
B	719	HIS	-	expression tag	UNP G1JZH9
B	720	HIS	-	expression tag	UNP G1JZH9
B	721	HIS	-	expression tag	UNP G1JZH9
B	722	HIS	-	expression tag	UNP G1JZH9
B	723	HIS	-	expression tag	UNP G1JZH9
C	1	MET	-	initiating methionine	UNP G1JZH9
C	2	ARG	-	expression tag	UNP G1JZH9
C	3	VAL	-	expression tag	UNP G1JZH9
C	4	LYS	-	expression tag	UNP G1JZH9
C	5	GLU	-	expression tag	UNP G1JZH9
C	6	LYS	-	expression tag	UNP G1JZH9
C	7	TYR	-	expression tag	UNP G1JZH9
C	8	GLN	-	expression tag	UNP G1JZH9
C	9	HIS	-	expression tag	UNP G1JZH9
C	10	LEU	-	expression tag	UNP G1JZH9
C	11	TRP	-	expression tag	UNP G1JZH9
C	12	ARG	-	expression tag	UNP G1JZH9
C	13	TRP	-	expression tag	UNP G1JZH9
C	14	GLY	-	expression tag	UNP G1JZH9
C	15	TRP	-	expression tag	UNP G1JZH9
C	16	ARG	-	expression tag	UNP G1JZH9
C	17	TRP	-	expression tag	UNP G1JZH9
C	18	GLY	-	expression tag	UNP G1JZH9
C	19	THR	-	expression tag	UNP G1JZH9
C	31	THR	VAL	conflict	UNP G1JZH9
C	33	LYS	ASN	conflict	UNP G1JZH9
C	114	GLU	GLN	conflict	UNP G1JZH9
C	533	VAL	ALA	conflict	UNP G1JZH9
C	536	MET	ILE	conflict	UNP G1JZH9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	544	GLN	LEU	conflict	UNP G1JZH9
C	568	LYS	GLN	conflict	UNP G1JZH9
C	583	THR	ALA	conflict	UNP G1JZH9
C	716	GLY	-	expression tag	UNP G1JZH9
C	717	GLY	-	expression tag	UNP G1JZH9
C	718	GLY	-	expression tag	UNP G1JZH9
C	719	HIS	-	expression tag	UNP G1JZH9
C	720	HIS	-	expression tag	UNP G1JZH9
C	721	HIS	-	expression tag	UNP G1JZH9
C	722	HIS	-	expression tag	UNP G1JZH9
C	723	HIS	-	expression tag	UNP G1JZH9
C	724	HIS	-	expression tag	UNP G1JZH9
D	2	MET	-	initiating methionine	UNP G1JZH9
D	3	ARG	-	expression tag	UNP G1JZH9
D	4	VAL	-	expression tag	UNP G1JZH9
D	5	LYS	-	expression tag	UNP G1JZH9
D	6	GLU	-	expression tag	UNP G1JZH9
D	7	LYS	-	expression tag	UNP G1JZH9
D	8	TYR	-	expression tag	UNP G1JZH9
D	9	GLN	-	expression tag	UNP G1JZH9
D	10	HIS	-	expression tag	UNP G1JZH9
D	11	LEU	-	expression tag	UNP G1JZH9
D	12	TRP	-	expression tag	UNP G1JZH9
D	13	ARG	-	expression tag	UNP G1JZH9
D	14	TRP	-	expression tag	UNP G1JZH9
D	15	GLY	-	expression tag	UNP G1JZH9
D	16	TRP	-	expression tag	UNP G1JZH9
D	17	ARG	-	expression tag	UNP G1JZH9
D	18	TRP	-	expression tag	UNP G1JZH9
D	19	GLY	-	expression tag	UNP G1JZH9
D	20	THR	-	expression tag	UNP G1JZH9
D	32	THR	VAL	conflict	UNP G1JZH9
D	34	LYS	ASN	conflict	UNP G1JZH9
D	115	GLU	GLN	conflict	UNP G1JZH9
D	532	VAL	ALA	conflict	UNP G1JZH9
D	535	MET	ILE	conflict	UNP G1JZH9
D	543	GLN	LEU	conflict	UNP G1JZH9
D	567	LYS	GLN	conflict	UNP G1JZH9
D	582	THR	ALA	conflict	UNP G1JZH9
D	715	GLY	-	expression tag	UNP G1JZH9
D	716	GLY	-	expression tag	UNP G1JZH9
D	717	GLY	-	expression tag	UNP G1JZH9

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Chain	Residue	Modelled	Actual	Comment	Reference
D	718	HIS	-	expression tag	UNP G1JZH9
D	719	HIS	-	expression tag	UNP G1JZH9
D	720	HIS	-	expression tag	UNP G1JZH9
D	721	HIS	-	expression tag	UNP G1JZH9
D	722	HIS	-	expression tag	UNP G1JZH9
D	723	HIS	-	expression tag	UNP G1JZH9
E	1	MET	-	initiating methionine	UNP G1JZH9
E	2	ARG	-	expression tag	UNP G1JZH9
E	3	VAL	-	expression tag	UNP G1JZH9
E	4	LYS	-	expression tag	UNP G1JZH9
E	5	GLU	-	expression tag	UNP G1JZH9
E	6	LYS	-	expression tag	UNP G1JZH9
E	7	TYR	-	expression tag	UNP G1JZH9
E	8	GLN	-	expression tag	UNP G1JZH9
E	9	HIS	-	expression tag	UNP G1JZH9
E	10	LEU	-	expression tag	UNP G1JZH9
E	11	TRP	-	expression tag	UNP G1JZH9
E	12	ARG	-	expression tag	UNP G1JZH9
E	13	TRP	-	expression tag	UNP G1JZH9
E	14	GLY	-	expression tag	UNP G1JZH9
E	15	TRP	-	expression tag	UNP G1JZH9
E	16	ARG	-	expression tag	UNP G1JZH9
E	17	TRP	-	expression tag	UNP G1JZH9
E	18	GLY	-	expression tag	UNP G1JZH9
E	19	THR	-	expression tag	UNP G1JZH9
E	31	THR	VAL	conflict	UNP G1JZH9
E	33	LYS	ASN	conflict	UNP G1JZH9
E	114	GLU	GLN	conflict	UNP G1JZH9
E	533	VAL	ALA	conflict	UNP G1JZH9
E	536	MET	ILE	conflict	UNP G1JZH9
E	544	GLN	LEU	conflict	UNP G1JZH9
E	568	LYS	GLN	conflict	UNP G1JZH9
E	583	THR	ALA	conflict	UNP G1JZH9
E	716	GLY	-	expression tag	UNP G1JZH9
E	717	GLY	-	expression tag	UNP G1JZH9
E	718	GLY	-	expression tag	UNP G1JZH9
E	719	HIS	-	expression tag	UNP G1JZH9
E	720	HIS	-	expression tag	UNP G1JZH9
E	721	HIS	-	expression tag	UNP G1JZH9
E	722	HIS	-	expression tag	UNP G1JZH9
E	723	HIS	-	expression tag	UNP G1JZH9
E	724	HIS	-	expression tag	UNP G1JZH9

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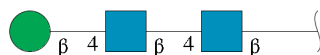
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP G1JZH9
A	2	ARG	-	expression tag	UNP G1JZH9
A	3	VAL	-	expression tag	UNP G1JZH9
A	4	LYS	-	expression tag	UNP G1JZH9
A	5	GLU	-	expression tag	UNP G1JZH9
A	6	LYS	-	expression tag	UNP G1JZH9
A	7	TYR	-	expression tag	UNP G1JZH9
A	8	GLN	-	expression tag	UNP G1JZH9
A	9	HIS	-	expression tag	UNP G1JZH9
A	10	LEU	-	expression tag	UNP G1JZH9
A	11	TRP	-	expression tag	UNP G1JZH9
A	12	ARG	-	expression tag	UNP G1JZH9
A	13	TRP	-	expression tag	UNP G1JZH9
A	14	GLY	-	expression tag	UNP G1JZH9
A	15	TRP	-	expression tag	UNP G1JZH9
A	16	ARG	-	expression tag	UNP G1JZH9
A	17	TRP	-	expression tag	UNP G1JZH9
A	18	GLY	-	expression tag	UNP G1JZH9
A	19	THR	-	expression tag	UNP G1JZH9
A	31	THR	VAL	conflict	UNP G1JZH9
A	33	LYS	ASN	conflict	UNP G1JZH9
A	114	GLU	GLN	conflict	UNP G1JZH9
A	533	VAL	ALA	conflict	UNP G1JZH9
A	536	MET	ILE	conflict	UNP G1JZH9
A	544	GLN	LEU	conflict	UNP G1JZH9
A	568	LYS	GLN	conflict	UNP G1JZH9
A	583	THR	ALA	conflict	UNP G1JZH9
A	716	GLY	-	expression tag	UNP G1JZH9
A	717	GLY	-	expression tag	UNP G1JZH9
A	718	GLY	-	expression tag	UNP G1JZH9
A	719	HIS	-	expression tag	UNP G1JZH9
A	720	HIS	-	expression tag	UNP G1JZH9
A	721	HIS	-	expression tag	UNP G1JZH9
A	722	HIS	-	expression tag	UNP G1JZH9
A	723	HIS	-	expression tag	UNP G1JZH9
A	724	HIS	-	expression tag	UNP G1JZH9
F	2	MET	-	initiating methionine	UNP G1JZH9
F	3	ARG	-	expression tag	UNP G1JZH9
F	4	VAL	-	expression tag	UNP G1JZH9
F	5	LYS	-	expression tag	UNP G1JZH9
F	6	GLU	-	expression tag	UNP G1JZH9
F	7	LYS	-	expression tag	UNP G1JZH9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	8	TYR	-	expression tag	UNP G1JZH9
F	9	GLN	-	expression tag	UNP G1JZH9
F	10	HIS	-	expression tag	UNP G1JZH9
F	11	LEU	-	expression tag	UNP G1JZH9
F	12	TRP	-	expression tag	UNP G1JZH9
F	13	ARG	-	expression tag	UNP G1JZH9
F	14	TRP	-	expression tag	UNP G1JZH9
F	15	GLY	-	expression tag	UNP G1JZH9
F	16	TRP	-	expression tag	UNP G1JZH9
F	17	ARG	-	expression tag	UNP G1JZH9
F	18	TRP	-	expression tag	UNP G1JZH9
F	19	GLY	-	expression tag	UNP G1JZH9
F	20	THR	-	expression tag	UNP G1JZH9
F	32	THR	VAL	conflict	UNP G1JZH9
F	34	LYS	ASN	conflict	UNP G1JZH9
F	115	GLU	GLN	conflict	UNP G1JZH9
F	532	VAL	ALA	conflict	UNP G1JZH9
F	535	MET	ILE	conflict	UNP G1JZH9
F	543	GLN	LEU	conflict	UNP G1JZH9
F	567	LYS	GLN	conflict	UNP G1JZH9
F	582	THR	ALA	conflict	UNP G1JZH9
F	715	GLY	-	expression tag	UNP G1JZH9
F	716	GLY	-	expression tag	UNP G1JZH9
F	717	GLY	-	expression tag	UNP G1JZH9
F	718	HIS	-	expression tag	UNP G1JZH9
F	719	HIS	-	expression tag	UNP G1JZH9
F	720	HIS	-	expression tag	UNP G1JZH9
F	721	HIS	-	expression tag	UNP G1JZH9
F	722	HIS	-	expression tag	UNP G1JZH9
F	723	HIS	-	expression tag	UNP G1JZH9

- Molecule 2 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				AltConf	Trace
2	G	3	Total	C	N	O	0	0
			39	22	2	15		

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Mol	Chain	Residues	Atoms				AltConf	Trace
2	I	3	Total	C	N	O	0	0
			39	22	2	15		
2	N	3	Total	C	N	O	0	0
			39	22	2	15		
2	R	3	Total	C	N	O	0	0
			39	22	2	15		
2	T	3	Total	C	N	O	0	0
			39	22	2	15		
2	Y	3	Total	C	N	O	0	0
			39	22	2	15		
2	Z	3	Total	C	N	O	0	0
			39	22	2	15		
2	c	3	Total	C	N	O	0	0
			39	22	2	15		
2	d	3	Total	C	N	O	0	0
			39	22	2	15		
2	g	3	Total	C	N	O	0	0
			39	22	2	15		
2	j	3	Total	C	N	O	0	0
			39	22	2	15		
2	k	3	Total	C	N	O	0	0
			39	22	2	15		
2	n	3	Total	C	N	O	0	0
			39	22	2	15		
2	o	3	Total	C	N	O	0	0
			39	22	2	15		
2	s	3	Total	C	N	O	0	0
			39	22	2	15		
2	t	3	Total	C	N	O	0	0
			39	22	2	15		
2	u	3	Total	C	N	O	0	0
			39	22	2	15		
2	x	3	Total	C	N	O	0	0
			39	22	2	15		
2	y	3	Total	C	N	O	0	0
			39	22	2	15		
2	z	3	Total	C	N	O	0	0
			39	22	2	15		
2	0	3	Total	C	N	O	0	0
			39	22	2	15		
2	1	3	Total	C	N	O	0	0
			39	22	2	15		

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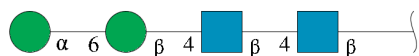
Mol	Chain	Residues	Atoms				AltConf	Trace
2	4	3	Total	C	N	O	0	0
			39	22	2	15		
2	8	3	Total	C	N	O	0	0
			39	22	2	15		
2	AA	3	Total	C	N	O	0	0
			39	22	2	15		
2	BA	3	Total	C	N	O	0	0
			39	22	2	15		

- 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				AltConf	Trace
3	H	2	Total	C	N	O	0	0
			28	16	2	10		
3	J	2	Total	C	N	O	0	0
			28	16	2	10		
3	M	2	Total	C	N	O	0	0
			28	16	2	10		
3	O	2	Total	C	N	O	0	0
			28	16	2	10		
3	P	2	Total	C	N	O	0	0
			28	16	2	10		
3	S	2	Total	C	N	O	0	0
			28	16	2	10		
3	b	2	Total	C	N	O	0	0
			28	16	2	10		
3	h	2	Total	C	N	O	0	0
			28	16	2	10		
3	v	2	Total	C	N	O	0	0
			28	16	2	10		
3	EA	2	Total	C	N	O	0	0
			28	16	2	10		

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



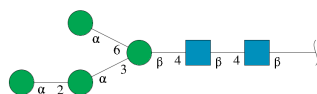
Mol	Chain	Residues	Atoms				AltConf	Trace
4	K	4	Total	C	N	O	0	0
			50	28	2	20		
4	i	4	Total	C	N	O	0	0
			50	28	2	20		
4	p	4	Total	C	N	O	0	0
			50	28	2	20		
4	5	4	Total	C	N	O	0	0
			50	28	2	20		

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



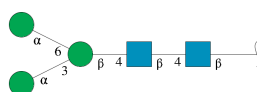
Mol	Chain	Residues	Atoms				AltConf	Trace
5	L	4	Total	C	N	O	0	0
			50	28	2	20		
5	U	4	Total	C	N	O	0	0
			50	28	2	20		
5	X	4	Total	C	N	O	0	0
			50	28	2	20		
5	m	4	Total	C	N	O	0	0
			50	28	2	20		
5	r	4	Total	C	N	O	0	0
			50	28	2	20		
5	3	4	Total	C	N	O	0	0
			50	28	2	20		
5	9	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-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.



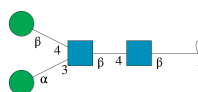
Mol	Chain	Residues	Atoms				AltConf	Trace
6	Q	6	Total	C	N	O	0	0
			72	40	2	30		
6	2	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 7 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				AltConf	Trace
7	V	5	Total	C	N	O	0	0
			61	34	2	25		
7	e	5	Total	C	N	O	0	0
			61	34	2	25		
7	f	5	Total	C	N	O	0	0
			61	34	2	25		
7	w	5	Total	C	N	O	0	0
			61	34	2	25		
7	6	5	Total	C	N	O	0	0
			61	34	2	25		

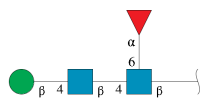
- Molecule 8 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[beta-D-mannopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
8	W	4	Total	C	N	O	0	0
			50	28	2	20		

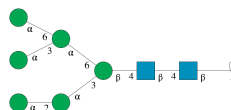
- Molecule 9 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-b

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



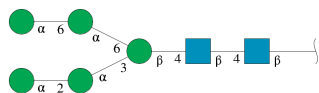
Mol	Chain	Residues	Atoms				AltConf	Trace
9	a	4	Total	C	N	O	0	0
			49	28	2	19		
9	CA	4	Total	C	N	O	0	0
			49	28	2	19		

- Molecule 10 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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.



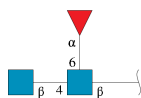
Mol	Chain	Residues	Atoms				AltConf	Trace
10	1	8	Total	C	N	O	0	0
			94	52	2	40		

- Molecule 11 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-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.



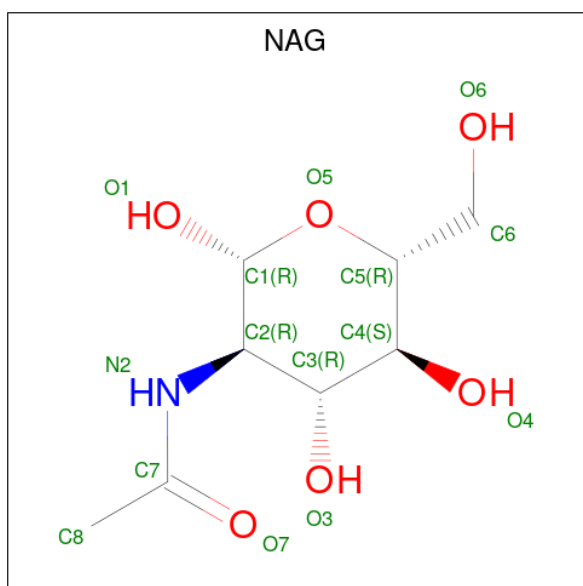
Mol	Chain	Residues	Atoms				AltConf	Trace
11	q	7	Total	C	N	O	0	0
			83	46	2	35		
11	7	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 12 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				AltConf	Trace
12	DA	3	Total	C	N	O	0	0
			38	22	2	14		
12	FA	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 13 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



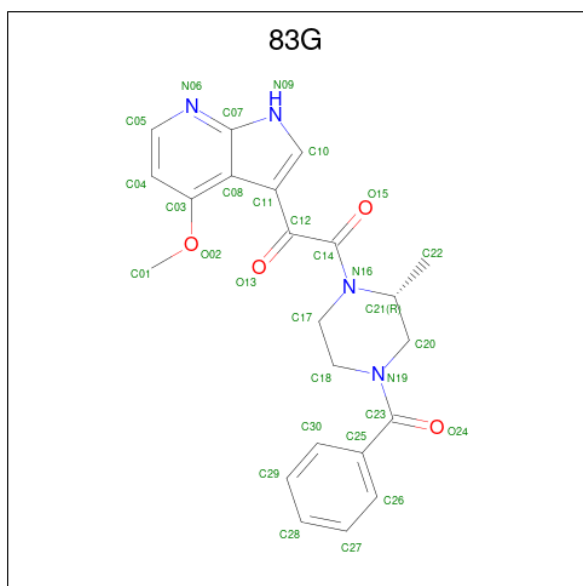
Mol	Chain	Residues	Atoms				AltConf
13	B	1	Total	C	N	O	0
			14	8	1	5	
13	C	1	Total	C	N	O	0
			14	8	1	5	
13	C	1	Total	C	N	O	0
			14	8	1	5	
13	C	1	Total	C	N	O	0
			14	8	1	5	
13	C	1	Total	C	N	O	0
			14	8	1	5	
13	D	1	Total	C	N	O	0
			14	8	1	5	
13	E	1	Total	C	N	O	0
			14	8	1	5	

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Mol	Chain	Residues	Atoms				AltConf
13	E	1	Total	C	N	O	0
			14	8	1	5	
13	A	1	Total	C	N	O	0
			14	8	1	5	
13	A	1	Total	C	N	O	0
			14	8	1	5	
13	A	1	Total	C	N	O	0
			14	8	1	5	

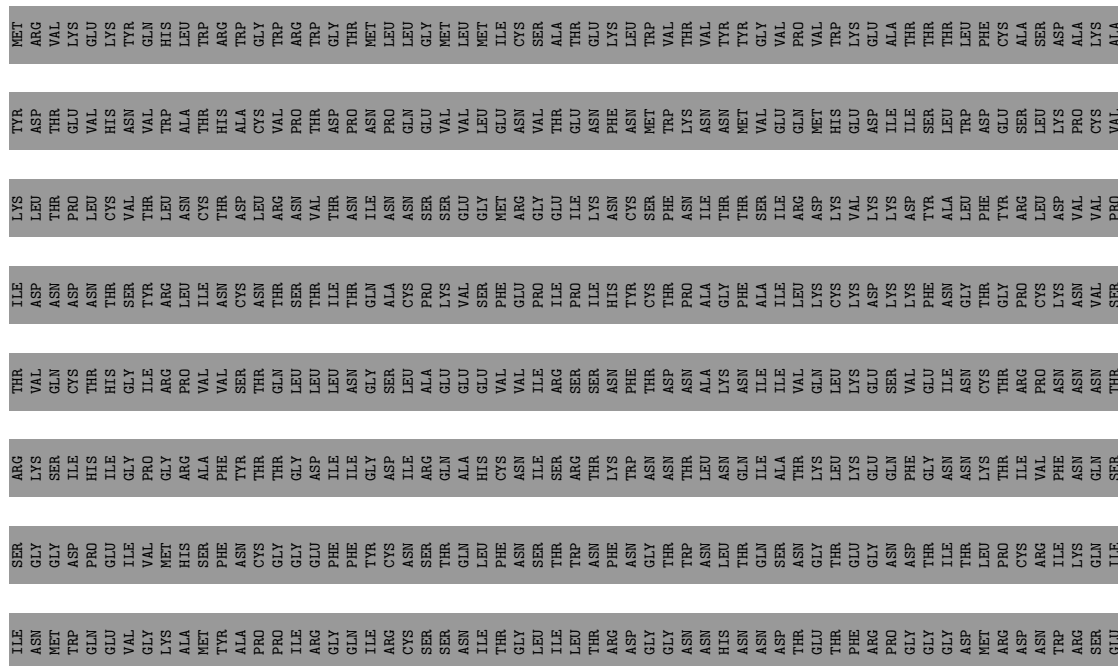
- Molecule 14 is 1-[(2R)-4-(benzenecarbonyl)-2-methylpiperazin-1-yl]-2-(4-methoxy-1H-pyrrolo[2,3-b]pyridin-3-yl)ethane-1,2-dione (CCD ID: 83G) (formula: C₂₂H₂₂N₄O₄).

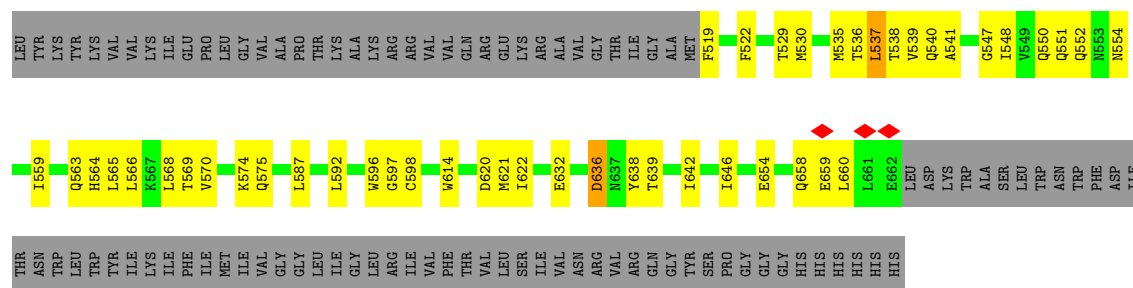


Mol	Chain	Residues	Atoms				AltConf
14	C	1	Total	C	N	O	0
			30	22	4	4	
14	E	1	Total	C	N	O	0
			30	22	4	4	
14	A	1	Total	C	N	O	0
			30	22	4	4	

- Molecule 1: Envelope glycoprotein gp160

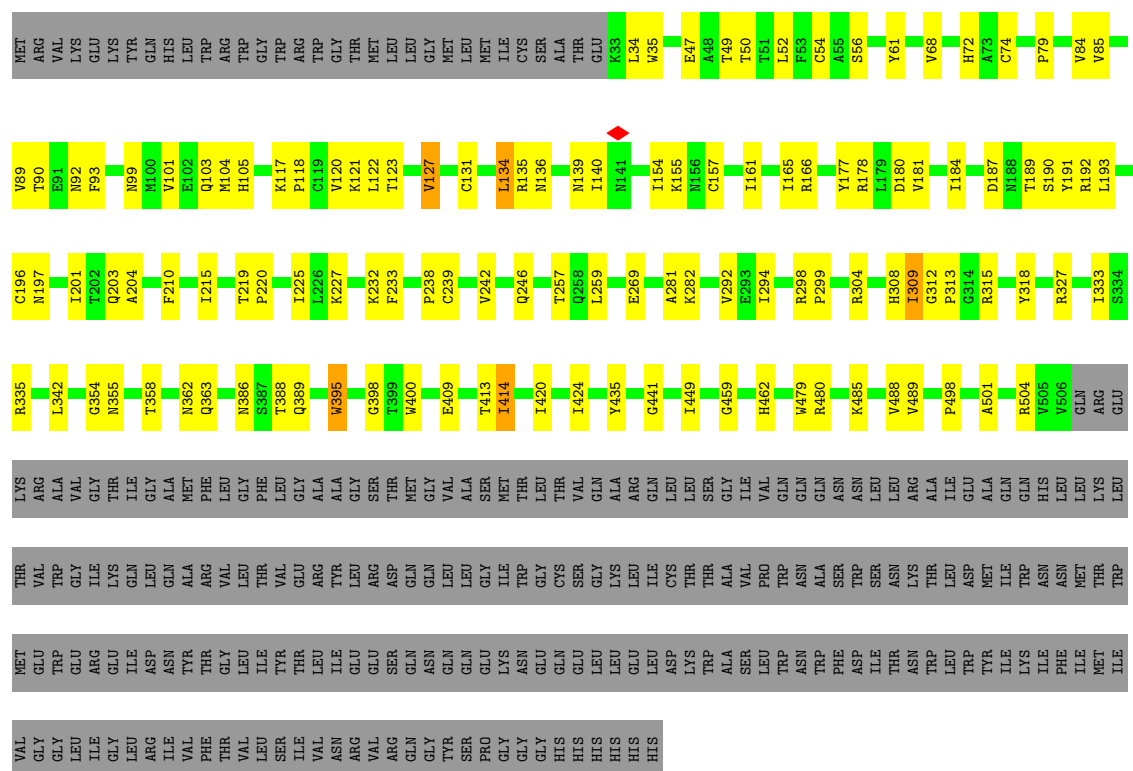






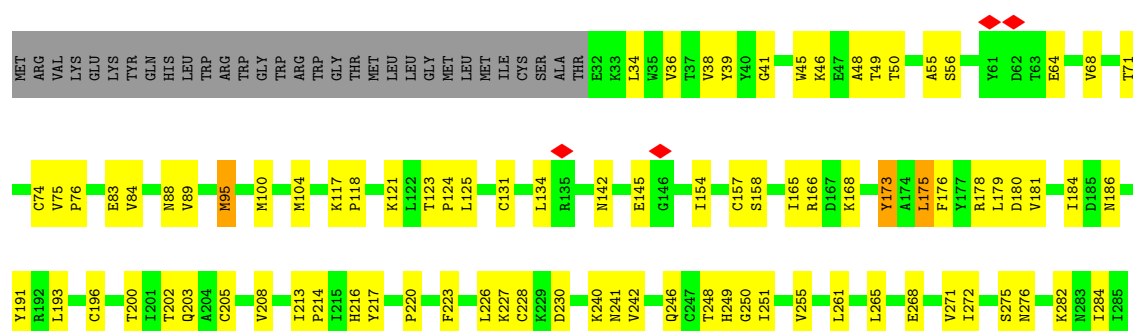
• Molecule 1: Envelope glycoprotein gp160

Chain E: 49% 16% 35%

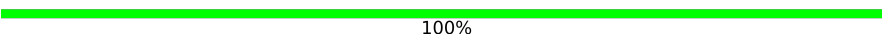


• Molecule 1: Envelope glycoprotein gp160

Chain A: 47% 18% 34%

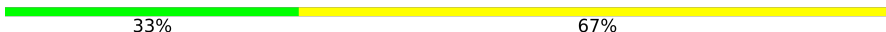


- Molecule 2: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  100%

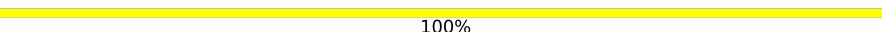


- Molecule 2: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  33% 67%



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

Chain N:  100%



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

Chain R:  67% 33%



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

Chain Y:  67% 33%



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

Chain Z:  67% 33%



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

Chain c:  100%



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

Chain d:  67% 33%



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

Chain g:  33% 67% 33%



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

Chain j:  33% 67%



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

Chain k:  100%



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

Chain n:  33% 100%



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

Chain o:  100%

MAG1
MAG2
BMA3

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

Chain s:  67% 33%

MAG1
MAG2
BMA3

- Molecule 2: 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%

MAG1
MAG2
BMA3

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

Chain u:  33% 67%

MAG1
MAG2
BMA3

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

Chain x:  67% 33%

MAG1
MAG2
BMA3

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

Chain y:  33% 33% 33%

MAG1
MAG2
BMA3

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

Chain z:  67% 33%



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

Chain 0:



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

Chain 1:



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

Chain 4:



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

Chain 8:



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

Chain AA:



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

Chain BA:



- 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



- 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



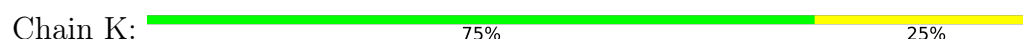
- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



- Molecule 4: 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



- Molecule 4: 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



- Molecule 4: 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:  100%



- Molecule 4: 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 5:  25% 25% 75%



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

Chain L:  50% 50%



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

Chain U:  50% 50%



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

Chain X:  25% 50% 50%



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

Chain m:  75% 25%



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

Chain r:  100%



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

Chain 3:  75% 25%



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

Chain 9:  50% 25% 25%



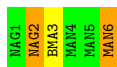
- Molecule 6: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  67% 17% 17%



- Molecule 6: alpha-D-mannopyranose-(1-2)-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 2:  50% 17% 33%

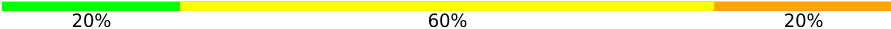


- Molecule 7: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain V:  20% 80%



- Molecule 7: 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 e: 



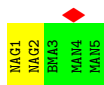
- Molecule 7: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain f: 



- Molecule 7: 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 w: 



- Molecule 7: 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 6: 



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

Chain W: 



- Molecule 9: 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 a: 



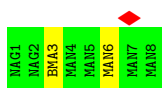
- Molecule 9: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain CA:  75% 25%



- Molecule 10: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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 l:  12% 75% 25%



- Molecule 11: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-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:  29% 71%



- Molecule 11: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-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 7:  14% 43% 57%



- Molecule 12: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain DA:  67% 100%



- Molecule 12: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain FA:  67% 33%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	50646	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	51.7	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.140	Depositor
Minimum map value	-0.052	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.02	Depositor
Map size (Å)	297.0, 297.0, 297.0	wwPDB
Map dimensions	180, 180, 180	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.65, 1.65, 1.65	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FUC, BMA, NAG, MAN, 83G

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	0.96	0/3859	1.11	4/5242 (0.1%)
1	B	0.89	0/1431	1.15	0/1947
1	C	0.94	0/3830	1.10	1/5205 (0.0%)
1	D	0.91	0/1179	1.18	1/1599 (0.1%)
1	E	0.93	0/3801	1.10	1/5167 (0.0%)
1	F	0.91	0/1179	1.19	0/1599
All	All	0.93	0/15279	1.12	7/20759 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	407	GLY	CA-C-O	-6.09	118.26	122.22
1	E	441	GLY	CA-C-O	-5.80	118.27	122.45
1	A	186	ASN	N-CA-C	-5.41	107.05	113.38
1	C	410	GLY	CA-C-O	-5.26	118.03	122.29
1	A	275	SER	N-CA-C	-5.25	105.24	110.97
1	A	145	GLU	CB-CA-C	-5.22	110.58	116.63
1	D	659	GLU	N-CA-C	-5.07	107.23	113.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	3780	0	3683	76	0
1	B	1399	0	1389	26	0
1	C	3751	0	3651	90	0
1	D	1159	0	1151	33	0
1	E	3722	0	3624	75	0
1	F	1159	0	1152	33	0
2	0	39	0	34	0	0
2	1	39	0	34	1	0
2	4	39	0	34	0	0
2	8	39	0	34	0	0
2	AA	39	0	34	0	0
2	BA	39	0	34	0	0
2	G	39	0	34	0	0
2	I	39	0	34	0	0
2	N	39	0	34	0	0
2	R	39	0	34	0	0
2	T	39	0	34	0	0
2	Y	39	0	34	0	0
2	Z	39	0	34	0	0
2	c	39	0	34	0	0
2	d	39	0	34	0	0
2	g	39	0	34	0	0
2	j	39	0	34	0	0
2	k	39	0	34	0	0
2	n	39	0	34	0	0
2	o	39	0	34	0	0
2	s	39	0	34	0	0
2	t	39	0	34	0	0
2	u	39	0	34	0	0
2	x	39	0	34	0	0
2	y	39	0	34	1	0
2	z	39	0	34	0	0
3	EA	28	0	25	0	0
3	H	28	0	25	0	0
3	J	28	0	25	0	0
3	M	28	0	25	0	0
3	O	28	0	25	1	0
3	P	28	0	25	1	0
3	S	28	0	25	0	0
3	b	28	0	25	0	0
3	h	28	0	25	0	0
3	v	28	0	25	0	0
4	5	50	0	43	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	K	50	0	43	0	0
4	i	50	0	43	0	0
4	p	50	0	43	0	0
5	3	50	0	43	0	0
5	9	50	0	43	2	0
5	L	50	0	43	0	0
5	U	50	0	43	0	0
5	X	50	0	43	0	0
5	m	50	0	43	2	0
5	r	50	0	43	0	0
6	2	72	0	61	2	0
6	Q	72	0	61	1	0
7	6	61	0	52	0	0
7	V	61	0	52	3	0
7	e	61	0	52	2	0
7	f	61	0	52	0	0
7	w	61	0	52	0	0
8	W	50	0	43	0	0
9	CA	49	0	43	0	0
9	a	49	0	43	0	0
10	l	94	0	79	1	0
11	7	83	0	70	4	0
11	q	83	0	70	6	0
12	DA	38	0	34	0	0
12	FA	38	0	34	0	0
13	A	42	0	39	0	0
13	B	14	0	13	0	0
13	C	56	0	52	0	0
13	D	14	0	13	0	0
13	E	28	0	26	0	0
14	A	30	0	0	0	0
14	C	30	0	0	0	0
14	E	30	0	0	0	0
All	All	17991	0	17198	325	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 9.

All (325) 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)
7:V:3:BMA:H2	7:V:4:MAN:H3	1.65	0.78
1:C:133:ASP:HB3	1:C:156:ASN:HB3	1.71	0.71
1:C:231:LYS:HE2	3:P:1:NAG:H5	1.73	0.70
1:C:227:LYS:HE3	1:C:485:LYS:HE3	1.73	0.69
1:C:129:LEU:HB3	1:C:157:CYS:SG	2.33	0.68
1:C:357:LYS:HA	1:C:464(A):ASP:HB3	1.75	0.67
1:E:184:ILE:HG13	1:E:192:ARG:HB2	1.77	0.66
7:e:3:BMA:H2	7:e:4:MAN:H3	1.78	0.66
1:B:542:ARG:HG2	1:A:41:GLY:HA3	1.79	0.63
1:F:557:ARG:HG2	1:F:559:ILE:H	1.63	0.63
1:C:134:LEU:HD13	1:C:152:GLY:HA3	1.80	0.62
6:2:3:BMA:H61	6:2:6:MAN:H3	1.81	0.62
1:E:193:LEU:HB2	1:E:196:CYS:HB2	1.83	0.61
1:B:583:VAL:HG13	1:D:587:LEU:HD11	1.83	0.61
1:C:117:LYS:HB3	1:C:118:PRO:HD3	1.83	0.60
1:A:48:ALA:HB3	1:A:490:LYS:HB3	1.84	0.60
1:E:187:ASP:HB2	1:E:190:SER:HB3	1.84	0.59
1:C:131:CYS:HA	1:C:157:CYS:HA	1.83	0.59
1:B:542:ARG:HD3	1:B:602:LEU:HD13	1.83	0.58
1:B:617:LYS:HB2	1:B:622:ILE:HG22	1.84	0.58
11:7:3:BMA:H3	11:7:4:MAN:H5	1.84	0.58
11:q:3:BMA:H3	11:q:4:MAN:H5	1.85	0.58
1:C:68:VAL:HG13	1:C:69:TRP:HD1	1.69	0.57
1:F:555:LEU:HD22	1:F:557:ARG:HH11	1.70	0.56
1:C:209:SER:OG	1:C:211:GLU:HG2	2.06	0.56
1:A:276:ASN:HB3	1:A:282:LYS:HG3	1.87	0.56
1:C:135:ARG:NH2	1:C:326:ILE:HG13	2.20	0.56
1:E:131:CYS:HA	1:E:157:CYS:HA	1.87	0.56
1:A:220:PRO:HG2	1:A:223:PHE:CD1	2.41	0.55
1:E:61:TYR:CZ	1:F:557:ARG:HD3	2.42	0.54
1:A:117:LYS:HB3	1:A:118:PRO:HD3	1.89	0.54
1:E:56:SER:H	1:E:74:CYS:HB3	1.71	0.54
1:A:121:LYS:HA	1:A:202:THR:HA	1.90	0.54
1:A:388:THR:HG21	5:9:1:NAG:H62	1.89	0.54
1:F:520:LEU:HD12	1:F:536:THR:HG21	1.90	0.54
1:E:299:PRO:HB2	1:E:327:ARG:HB2	1.90	0.53
1:F:562:GLN:O	1:F:563:GLN:C	2.50	0.53
1:A:230:ASP:HA	1:A:241:ASN:HB2	1.90	0.53
1:E:203:GLN:HA	1:E:435:TYR:HB3	1.90	0.53
1:B:535:MET:HA	1:B:538:THR:HG23	1.89	0.53
1:E:187:ASP:HB2	1:E:190:SER:CB	2.39	0.53
1:B:527:GLY:HA3	1:A:88:ASN:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:69:TRP:HZ3	1:C:108:ILE:HG23	1.73	0.53
1:E:72:HIS:O	1:F:557:ARG:NH2	2.41	0.53
1:C:137:VAL:C	1:C:139:ASN:H	2.16	0.53
1:A:381:GLU:HG2	1:A:439:ILE:HG12	1.90	0.53
1:A:249:HIS:O	1:A:251:ILE:HG13	2.09	0.52
1:B:521:GLY:HA2	1:B:548:ILE:HG21	1.90	0.52
1:C:254:VAL:HG13	6:Q:1:NAG:H81	1.91	0.52
1:C:135:ARG:NE	1:C:150:MET:HB3	2.25	0.52
1:E:304:ARG:HD3	1:E:318:TYR:CD2	2.43	0.52
1:B:565:LEU:C	1:B:567:LYS:H	2.18	0.51
1:C:331:CYS:HB2	1:C:416:LEU:HB2	1.92	0.51
1:C:377:ASN:OD1	1:C:381:GLU:C	2.53	0.51
1:F:534:SER:HA	1:F:537:LEU:HD13	1.90	0.51
1:E:333:ILE:HB	1:E:414:ILE:HB	1.91	0.51
1:D:642:ILE:O	1:D:646:ILE:HG12	2.09	0.51
1:C:343:ASN:HD21	1:C:400:TRP:CD1	2.29	0.51
1:B:529:THR:HA	1:B:627:THR:HA	1.93	0.51
1:A:45:TRP:HB2	1:A:490:LYS:O	2.11	0.51
1:D:550:GLN:HG2	1:F:588:ARG:HG3	1.93	0.50
1:C:295:ASN:HB2	1:C:332:ASN:HB2	1.93	0.50
1:A:454:LEU:HD23	1:A:470:PRO:HA	1.94	0.50
1:E:34:LEU:HD13	1:E:498:PRO:HB2	1.94	0.50
1:C:334:SER:HB3	1:C:337:LYS:HB3	1.93	0.50
1:C:385:CYS:HA	1:C:418:CYS:HA	1.93	0.50
1:C:362:ASN:HB3	1:C:469:ARG:HG2	1.94	0.50
1:A:261:LEU:HB3	1:A:447:SER:HB2	1.94	0.50
1:F:555:LEU:C	1:F:557:ARG:H	2.19	0.50
1:A:95:MET:HG2	1:A:484:TYR:HA	1.94	0.50
1:C:72:HIS:HB3	1:D:565:LEU:HD13	1.93	0.49
1:E:333:ILE:O	1:E:414:ILE:N	2.45	0.49
1:D:550:GLN:NE2	1:D:554:ASN:HB2	2.26	0.49
1:E:135:ARG:H	1:E:154:ILE:HG21	1.78	0.49
1:A:39:TYR:HE2	1:A:497:ALA:HB3	1.75	0.49
1:A:71:THR:HA	1:A:74:CYS:CB	2.42	0.49
1:C:174:ALA:HB1	1:C:319:THR:OG1	2.11	0.49
1:E:196:CYS:SG	1:E:197:ASN:N	2.86	0.49
1:F:611:ASN:HB3	1:F:614:TRP:CD2	2.48	0.49
1:D:654:GLU:O	1:D:658:GLN:N	2.42	0.49
1:A:49:THR:H	1:F:554:ASN:HB3	1.78	0.49
1:C:335:ARG:HA	1:C:414:ILE:HD11	1.94	0.48
1:E:227:LYS:HA	1:E:485:LYS:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:133:ASP:HB3	1:C:156:ASN:CB	2.42	0.48
1:C:156:ASN:OD1	1:C:173:TYR:HB2	2.14	0.48
1:E:93:PHE:HB2	1:E:233:PHE:HZ	1.78	0.48
1:E:232:LYS:HE2	1:E:269:GLU:HG2	1.95	0.48
5:m:3:BMA:H3	5:m:4:MAN:H5	1.95	0.48
1:C:54:CYS:HB2	1:C:215:ILE:HG23	1.95	0.48
1:B:643:TYR:HA	1:B:647:GLU:HB3	1.95	0.48
1:C:294:ILE:HG12	1:C:296:CYS:SG	2.54	0.48
1:C:133:ASP:O	1:C:134:LEU:C	2.55	0.48
1:E:165:ILE:HD12	1:E:165:ILE:H	1.79	0.48
1:F:632:GLU:HA	1:F:635:ILE:HG22	1.95	0.48
1:D:596:TRP:O	1:D:597:GLY:C	2.56	0.47
1:E:177:TYR:O	1:E:178:ARG:C	2.57	0.47
1:E:292:VAL:HB	1:E:449:ILE:HB	1.95	0.47
1:A:227:LYS:HE3	1:A:485:LYS:HD2	1.96	0.47
1:C:258:GLN:NE2	1:C:371:ILE:O	2.45	0.47
1:C:350:LYS:HB3	1:C:355:ASN:HA	1.96	0.47
1:B:538:THR:O	1:B:539:VAL:C	2.57	0.47
1:D:564:HIS:CE1	1:E:50:THR:HA	2.49	0.47
1:A:71:THR:HA	1:A:74:CYS:HB3	1.95	0.47
1:A:369:PRO:HA	1:A:372:VAL:HG22	1.95	0.47
1:E:177:TYR:O	1:E:180:ASP:N	2.40	0.47
1:B:546:SER:O	1:B:550:GLN:NE2	2.47	0.47
1:C:101:VAL:HG11	1:C:480:ARG:HG3	1.95	0.47
1:D:519:PHE:HB2	1:D:522:PHE:HB3	1.97	0.47
1:A:123:THR:N	1:A:124:PRO:HD2	2.30	0.47
1:A:291:SER:HB2	1:A:448:ASN:HB3	1.97	0.47
1:B:622:ILE:HD13	1:A:498:PRO:HD2	1.97	0.47
1:E:161:ILE:HG12	1:E:309:ILE:HD13	1.96	0.47
1:F:570:VAL:O	1:F:574:LYS:HG2	2.15	0.47
1:F:555:LEU:HD22	1:F:557:ARG:NH1	2.29	0.47
5:m:3:BMA:H2	5:m:4:MAN:H3	1.96	0.47
1:D:570:VAL:O	1:D:574:LYS:HD3	2.15	0.47
1:A:131:CYS:SG	1:A:191:TYR:HB2	2.55	0.47
1:A:304:ARG:HH21	1:A:440:ARG:HG2	1.80	0.47
1:C:259:LEU:HB2	1:C:374:HIS:CE1	2.50	0.46
1:C:345:ILE:HD13	1:C:452:LEU:HD12	1.97	0.46
1:F:559:ILE:HA	1:F:563:GLN:HB2	1.96	0.46
1:C:121:LYS:HA	1:C:202:THR:HA	1.97	0.46
1:D:614:TRP:HA	1:D:638:TYR:CD2	2.50	0.46
1:E:413:THR:O	1:E:414:ILE:C	2.58	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:569:THR:O	1:D:570:VAL:C	2.58	0.46
1:D:638:TYR:O	1:D:639:THR:C	2.58	0.46
1:A:246:GLN:OE1	1:A:246:GLN:N	2.46	0.46
1:A:131:CYS:HA	1:A:157:CYS:HA	1.98	0.46
1:E:395:TRP:HB3	1:E:400:TRP:HH2	1.80	0.46
1:E:498:PRO:HB3	1:F:610:TRP:CD2	2.50	0.46
1:A:319:THR:C	1:A:438:PRO:HG2	2.41	0.46
1:D:566:LEU:HB3	1:D:575:GLN:NE2	2.30	0.46
1:E:104:MET:O	1:E:105:HIS:C	2.57	0.46
1:A:205:CYS:O	1:A:208:VAL:HG22	2.15	0.46
1:C:344:GLN:O	1:C:348:LYS:HG2	2.16	0.46
1:E:239:CYS:HB3	1:E:242:VAL:HG22	1.98	0.46
1:C:56:SER:H	1:C:74:CYS:HB3	1.81	0.46
1:F:585:ARG:HG3	1:F:588:ARG:HH21	1.80	0.46
1:A:217:TYR:N	1:A:248:THR:OG1	2.49	0.45
7:V:3:BMA:H4	7:V:5:MAN:H3	1.99	0.45
1:B:630:GLU:HA	1:B:633:ARG:HD2	1.98	0.45
1:C:284:ILE:HB	1:C:454:LEU:HB2	1.97	0.45
1:C:292:VAL:HB	1:C:449:ILE:HB	1.97	0.45
1:A:249:HIS:O	1:A:251:ILE:N	2.49	0.45
1:C:68:VAL:HG13	1:C:69:TRP:CD1	2.51	0.45
1:C:349:LEU:HB2	1:C:359:ILE:HG12	1.99	0.45
1:C:145:GLU:C	1:C:150:MET:H	2.24	0.45
1:E:54:CYS:HB2	1:E:215:ILE:HG23	1.98	0.45
1:E:84:VAL:HG11	1:F:521:GLY:HA3	1.99	0.45
1:C:308:HIS:HA	1:C:316:ALA:HA	1.99	0.45
1:C:360:VAL:HG13	1:C:394:THR:HG22	1.99	0.45
1:E:52:LEU:N	1:E:103:GLN:OE1	2.49	0.45
1:D:564:HIS:HE2	1:E:49:THR:C	2.25	0.45
1:C:448:ASN:O	1:C:450:THR:HG23	2.16	0.45
1:E:90:THR:HA	1:E:239:CYS:O	2.17	0.45
1:A:181:VAL:HB	1:A:191:TYR:HB3	1.99	0.45
1:A:255:VAL:HG22	1:A:479:TRP:CH2	2.52	0.45
1:E:131:CYS:N	1:E:189:THR:O	2.42	0.45
1:F:614:TRP:CH2	1:F:642:ILE:HG12	2.51	0.45
1:E:35:TRP:HB2	1:E:501:ALA:O	2.17	0.45
1:A:387:SER:HA	1:A:390:LEU:HD23	1.99	0.45
7:e:3:BMA:H3	7:e:4:MAN:H5	1.99	0.45
1:B:530:MET:HE2	1:B:623:TRP:CD1	2.52	0.44
1:C:128:THR:OG1	1:A:168:LYS:NZ	2.48	0.44
1:C:221:ALA:HB2	1:D:552:GLN:HE22	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:ILE:HG12	1:C:488:VAL:HG12	1.99	0.44
1:E:134:LEU:HD23	1:E:136:ASN:HB2	1.98	0.44
1:E:398:GLY:HA2	11:q:6:MAN:H2	1.98	0.44
1:A:55:ALA:O	1:A:216:HIS:HB2	2.17	0.44
1:A:297:THR:HA	1:A:444:ARG:HA	1.98	0.44
1:B:618:THR:O	1:B:622:ILE:HG23	2.17	0.44
1:C:358:THR:HG23	1:C:395(A):ASN:HA	1.99	0.44
1:E:342:LEU:HB2	1:E:395:TRP:CH2	2.51	0.44
1:A:104:MET:HG2	1:A:479:TRP:HB3	1.99	0.44
1:E:204:ALA:HB1	1:E:210:PHE:HZ	1.82	0.44
1:E:386:ASN:C	1:E:388:THR:H	2.26	0.44
1:A:46:LYS:O	1:A:490:LYS:HG2	2.17	0.44
1:F:630:GLU:O	1:F:633:ARG:HG2	2.18	0.44
1:C:137:VAL:C	1:C:139:ASN:N	2.76	0.44
1:C:137:VAL:HG11	1:C:140:ILE:HG12	2.00	0.44
1:F:598:CYS:HB2	1:F:601:LYS:HB2	1.99	0.44
1:E:155:LYS:HE2	1:E:191:TYR:CE1	2.52	0.44
1:A:268:GLU:O	1:A:289:LYS:HD2	2.18	0.44
1:A:379:GLY:HA2	6:2:2:NAG:H62	2.00	0.44
1:B:550:GLN:N	1:B:550:GLN:HE21	2.16	0.44
1:C:353:PHE:HB3	1:C:357:LYS:HD2	2.00	0.44
1:E:47:GLU:HA	1:E:489:VAL:HG12	2.00	0.44
1:E:358:THR:HG23	1:E:395:TRP:O	2.16	0.44
1:C:231:LYS:O	1:C:267:GLU:HB3	2.18	0.44
1:E:61:TYR:CE2	1:F:557:ARG:HB2	2.53	0.44
1:C:184:ILE:C	1:C:186:ASN:H	2.26	0.44
1:E:354:GLY:O	1:E:355:ASN:C	2.60	0.43
1:B:604:CYS:SG	1:B:604:CYS:O	2.75	0.43
1:C:42:VAL:HG22	1:C:493:PRO:O	2.17	0.43
1:C:353:PHE:HE2	1:C:468:PHE:HZ	1.66	0.43
1:D:535:MET:O	1:D:538:THR:HG22	2.18	0.43
1:E:257:THR:C	1:E:259:LEU:H	2.26	0.43
1:F:632:GLU:O	1:F:633:ARG:C	2.61	0.43
11:7:3:BMA:H3	11:7:4:MAN:C5	2.47	0.43
1:B:565:LEU:C	1:B:567:LYS:N	2.76	0.43
1:E:101:VAL:HG11	1:E:480:ARG:HG3	2.00	0.43
1:E:117:LYS:HB3	1:E:118:PRO:HD3	1.99	0.43
10:l:3:BMA:H61	10:l:6:MAN:H2	1.83	0.43
1:C:225:ILE:HB	1:C:245:VAL:HG23	2.00	0.43
1:A:304:ARG:HB2	1:A:440:ARG:HB3	2.00	0.43
1:B:654:GLU:O	1:B:658:GLN:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:276:ASN:HB3	1:C:282:LYS:HG3	2.01	0.43
1:C:277:PHE:CE2	3:O:1:NAG:H81	2.53	0.43
1:A:178:ARG:O	1:A:180:ASP:N	2.42	0.43
1:B:532:VAL:O	1:B:535:MET:N	2.51	0.43
1:E:220:PRO:HB3	1:F:578:ALA:HB1	1.99	0.43
1:A:158:SER:HA	1:A:173:TYR:HA	2.00	0.43
7:V:3:BMA:O5	7:V:5:MAN:H2	2.19	0.43
11:q:3:BMA:H3	11:q:4:MAN:C5	2.47	0.43
2:1:1:NAG:H61	2:1:2:NAG:N2	2.32	0.43
1:D:566:LEU:HD22	1:D:566:LEU:H	1.84	0.43
1:A:75:VAL:HB	1:A:76:PRO:HD2	2.00	0.43
1:E:136:ASN:HB3	1:E:140:ILE:HB	2.01	0.43
1:A:176:PHE:HE2	1:A:317:PHE:HZ	1.66	0.43
1:A:203:GLN:HA	1:A:435:TYR:HB3	2.01	0.43
1:C:63:THR:HB	1:C:66:HIS:HD2	1.84	0.43
1:D:536:THR:O	1:D:537:LEU:C	2.62	0.43
1:A:49:THR:HB	1:F:560:GLU:HB3	1.99	0.43
1:D:563:GLN:HG3	1:D:564:HIS:N	2.34	0.43
1:D:632:GLU:O	1:D:636:ASP:N	2.51	0.43
1:E:120:VAL:HB	1:E:203:GLN:HG2	2.01	0.43
1:A:226:LEU:O	1:A:486:TYR:HA	2.19	0.43
1:A:342:LEU:HB3	1:A:395:TRP:CH2	2.54	0.43
1:C:153:GLU:O	1:C:178:ARG:N	2.50	0.42
1:C:61:TYR:CE1	1:D:565:LEU:HD11	2.54	0.42
1:C:261:LEU:O	1:C:262:ASN:C	2.62	0.42
1:A:175:LEU:HD23	1:A:175:LEU:HA	1.88	0.42
1:A:386:ASN:HB3	1:A:417:PRO:HD2	2.01	0.42
1:C:193:LEU:HD13	1:C:423:ILE:HD11	2.02	0.42
1:D:592:LEU:HB3	1:D:596:TRP:CZ2	2.55	0.42
1:E:362:ASN:OD1	1:E:363:GLN:N	2.52	0.42
1:A:50:THR:HA	1:F:561:ALA:CB	2.48	0.42
1:F:522:PHE:CE1	1:F:540:GLN:HA	2.54	0.42
1:C:307:ILE:O	1:C:317:PHE:N	2.48	0.42
1:C:476:ARG:HA	1:C:479:TRP:CD1	2.55	0.42
1:E:335:ARG:HB3	11:q:1:NAG:H81	2.00	0.42
11:q:4:MAN:H2	11:q:5:MAN:H2	1.80	0.42
1:B:604:CYS:SG	1:A:38:VAL:HB	2.60	0.42
1:C:104:MET:O	1:C:108:ILE:HG12	2.20	0.42
1:C:305:LYS:HB2	1:C:321:GLY:HA3	2.01	0.42
1:E:49:THR:HB	1:E:99:ASN:ND2	2.34	0.42
1:A:272:ILE:HG12	1:A:286:VAL:HG22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:527:GLY:HA3	1:A:88:ASN:CB	2.49	0.42
1:D:547:GLY:O	1:D:551:GLN:HG3	2.19	0.42
1:D:559:ILE:HG22	1:D:563:GLN:HE21	1.84	0.42
1:E:184:ILE:CG1	1:E:192:ARG:HB2	2.49	0.42
1:C:69:TRP:CZ3	1:C:108:ILE:HD12	2.55	0.42
1:E:312:GLY:HA3	1:E:315:ARG:HD2	2.01	0.42
1:C:96:TRP:C	1:C:98:ASN:H	2.27	0.42
1:C:281:ALA:O	1:C:282:LYS:C	2.63	0.42
1:C:298:ARG:NH1	1:C:327:ARG:O	2.53	0.42
1:D:596:TRP:HD1	1:D:598:CYS:HB2	1.84	0.42
1:E:281:ALA:O	1:E:282:LYS:C	2.62	0.42
1:A:178:ARG:HG2	1:A:179:LEU:H	1.85	0.42
4:5:3:BMA:H62	4:5:4:MAN:H2	1.86	0.42
1:C:58:ALA:C	1:C:60:ALA:N	2.77	0.42
1:E:135:ARG:N	1:E:154:ILE:HG21	2.34	0.42
1:E:308:HIS:NE2	1:E:312:GLY:O	2.53	0.42
1:E:389:GLN:HB2	1:E:414:ILE:CG2	2.49	0.42
1:A:362:ASN:HB3	1:A:469:ARG:HG2	2.01	0.42
1:A:388:THR:HG21	5:9:1:NAG:C6	2.50	0.42
1:B:663:LEU:H	1:B:663:LEU:HG	1.77	0.42
1:C:369:PRO:O	1:C:370:GLU:C	2.63	0.42
1:C:407:GLY:O	1:C:408:THR:C	2.63	0.42
1:E:166:ARG:HH22	1:E:313:PRO:HD3	1.85	0.42
1:A:165:ILE:HD12	1:A:168:LYS:HB2	2.02	0.42
1:A:292:VAL:HB	1:A:449:ILE:HB	2.02	0.42
1:C:73:ALA:O	1:C:74:CYS:C	2.63	0.41
1:C:119:CYS:HB3	1:C:203:GLN:O	2.20	0.41
1:E:225:ILE:HG12	1:E:488:VAL:HG22	2.02	0.41
1:A:425:ASN:HA	1:A:433:ALA:HA	2.02	0.41
1:C:142:ASN:HD22	1:C:143:SER:N	2.19	0.41
1:A:100:MET:SD	1:A:488:VAL:HG23	2.61	0.41
11:7:4:MAN:H2	11:7:5:MAN:H2	1.79	0.41
1:C:321(A):ASP:OD1	1:C:322:ILE:N	2.52	0.41
1:C:509:GLU:HB2	1:D:660:LEU:HD12	2.02	0.41
1:E:177:TYR:CD2	1:E:420:ILE:HB	2.56	0.41
1:A:142:ASN:N	1:A:142:ASN:OD1	2.53	0.41
1:C:272:ILE:HD12	1:C:286:VAL:HG22	2.01	0.41
1:A:45:TRP:HH2	1:A:89:VAL:HG11	1.86	0.41
11:q:1:NAG:H61	11:q:2:NAG:N2	2.36	0.41
2:y:1:NAG:H62	2:y:2:NAG:O7	2.21	0.41
1:C:123:THR:HG23	1:A:166:ARG:HH21	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:424:ILE:HD12	1:E:424:ILE:HA	1.98	0.41
1:F:642:ILE:O	1:F:646:ILE:HG13	2.20	0.41
1:D:548:ILE:HD12	1:D:551:GLN:HB2	2.02	0.41
1:E:101:VAL:HG13	1:E:479:TRP:HB2	2.03	0.41
1:E:120:VAL:HG12	1:E:122:LEU:H	1.85	0.41
1:E:342:LEU:HB2	1:E:395:TRP:HH2	1.86	0.41
1:A:363:GLN:HG2	1:A:388:THR:HG23	2.02	0.41
1:A:213:ILE:O	1:A:214:PRO:C	2.63	0.41
1:B:656:ASN:ND2	1:F:535:MET:HG2	2.36	0.41
1:C:41:GLY:N	1:C:493:PRO:O	2.54	0.41
1:D:568:LEU:HD21	1:F:568:LEU:HD13	2.03	0.41
1:E:121:LYS:C	1:E:123:THR:N	2.78	0.41
1:E:459:GLY:O	1:E:462:HIS:ND1	2.54	0.41
1:A:272:ILE:CD1	1:A:349:LEU:HG	2.51	0.41
1:A:272:ILE:HG23	1:A:284:ILE:HG22	2.03	0.41
1:F:657:GLU:O	1:F:661:LEU:HG	2.21	0.41
11:7:1:NAG:H61	11:7:2:NAG:N2	2.36	0.41
1:C:132:THR:O	1:C:133:ASP:C	2.64	0.41
1:E:294:ILE:HB	1:E:333:ILE:HD11	2.02	0.41
1:F:655:LYS:O	1:F:659:GLU:HG3	2.21	0.41
1:C:288:LEU:HD21	1:C:345:ILE:HD11	2.02	0.40
1:D:540:GLN:O	1:D:541:ALA:C	2.65	0.40
1:D:621:MET:O	1:D:622:ILE:C	2.63	0.40
1:E:134:LEU:HA	1:E:154:ILE:HB	2.02	0.40
1:A:56:SER:N	1:A:74:CYS:SG	2.94	0.40
1:A:176:PHE:CE2	1:A:317:PHE:HZ	2.39	0.40
1:C:285:ILE:HD13	1:C:285:ILE:HA	1.93	0.40
1:C:317:PHE:CE2	1:C:319:THR:HB	2.56	0.40
1:B:619:LEU:HD23	1:B:619:LEU:HA	1.83	0.40
1:C:114:GLU:CD	1:D:570:VAL:HG13	2.46	0.40
1:D:529:THR:O	1:D:530:MET:C	2.64	0.40
1:E:92:ASN:HA	1:E:238:PRO:HA	2.02	0.40
1:C:58:ALA:C	1:C:60:ALA:H	2.29	0.40
1:F:557:ARG:NE	1:F:559:ILE:HB	2.37	0.40
1:E:504:ARG:HH12	1:A:509:GLU:H	1.70	0.40
1:A:95:MET:CG	1:A:484:TYR:HA	2.51	0.40
1:A:125:LEU:HD21	1:A:176:PHE:HZ	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	476/722 (66%)	442 (93%)	32 (7%)	2 (0%)	30	62
1	B	168/722 (23%)	154 (92%)	13 (8%)	1 (1%)	21	54
1	C	473/722 (66%)	420 (89%)	48 (10%)	5 (1%)	11	43
1	D	142/722 (20%)	125 (88%)	17 (12%)	0	100	100
1	E	470/722 (65%)	424 (90%)	42 (9%)	4 (1%)	14	47
1	F	142/722 (20%)	130 (92%)	10 (7%)	2 (1%)	9	38
All	All	1871/4332 (43%)	1695 (91%)	162 (9%)	14 (1%)	20	51

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	141	ASN
1	E	127	VAL
1	A	250	GLY
1	B	568	LEU
1	C	143	SER
1	C	240	LYS
1	A	228	CYS
1	F	563	GLN
1	C	466	GLU
1	E	395	TRP
1	C	414	ILE
1	F	614	TRP
1	E	79	PRO
1	E	414	ILE

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	430/639 (67%)	400 (93%)	30 (7%)	14	39
1	B	150/639 (24%)	140 (93%)	10 (7%)	15	41
1	C	427/639 (67%)	406 (95%)	21 (5%)	22	48
1	D	125/639 (20%)	121 (97%)	4 (3%)	34	59
1	E	424/639 (66%)	411 (97%)	13 (3%)	35	59
1	F	125/639 (20%)	116 (93%)	9 (7%)	13	38
All	All	1681/3834 (44%)	1594 (95%)	87 (5%)	22	47

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	523	LEU
1	B	550	GLN
1	B	585	ARG
1	B	622	ILE
1	B	624	ASN
1	B	639	THR
1	B	644	THR
1	B	652	GLN
1	B	675	ILE
1	B	677	ASN
1	C	96	TRP
1	C	123	THR
1	C	132	THR
1	C	142	ASN
1	C	159	PHE
1	C	160	ASN
1	C	169	VAL
1	C	189	THR
1	C	196	CYS
1	C	200	THR
1	C	201	ILE
1	C	232	LYS
1	C	278	THR
1	C	402	LEU
1	C	403	THR
1	C	426	MET
1	C	467	THR

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Mol	Chain	Res	Type
1	C	475	MET
1	C	485	LYS
1	C	488	VAL
1	C	496	VAL
1	D	537	LEU
1	D	539	VAL
1	D	620	ASP
1	D	636	ASP
1	E	68	VAL
1	E	85	VAL
1	E	89	VAL
1	E	127	VAL
1	E	134	LEU
1	E	139	ASN
1	E	181	VAL
1	E	201	ILE
1	E	219	THR
1	E	246	GLN
1	E	298	ARG
1	E	309	ILE
1	E	409	GLU
1	A	34	LEU
1	A	36	VAL
1	A	64	GLU
1	A	68	VAL
1	A	83	GLU
1	A	84	VAL
1	A	95	MET
1	A	134	LEU
1	A	154	ILE
1	A	173	TYR
1	A	175	LEU
1	A	184	ILE
1	A	193	LEU
1	A	196	CYS
1	A	200	THR
1	A	240	LYS
1	A	242	VAL
1	A	265	LEU
1	A	271	VAL
1	A	289	LYS
1	A	326	ILE

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Mol	Chain	Res	Type
1	A	333	ILE
1	A	360	VAL
1	A	373	MET
1	A	390	LEU
1	A	396	PHE
1	A	442	GLN
1	A	447	SER
1	A	461	ASN
1	A	500	LYS
1	F	545	LEU
1	F	564	HIS
1	F	569	THR
1	F	599	SER
1	F	626	MET
1	F	632	GLU
1	F	638	TYR
1	F	639	THR
1	F	647	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	543	GLN
1	B	671	ASN
1	B	677	ASN
1	C	66	HIS
1	C	72	HIS
1	C	99	ASN
1	C	249	HIS
1	C	395(A)	ASN
1	C	404	GLN
1	D	552	GLN
1	D	651	ASN
1	E	67	ASN
1	E	94	ASN
1	E	397	ASN
1	A	99	ASN
1	A	330	HIS
1	A	352	GLN
1	A	355	ASN
1	F	562	GLN
1	F	575	GLN

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Mol	Chain	Res	Type
1	F	591	GLN
1	F	624	ASN
1	F	652	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 ⓘ

219 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	0	1	1,2	14,14,15	0.38	0	17,19,21	0.71	0
2	NAG	0	2	2	14,14,15	0.39	0	17,19,21	0.96	1 (5%)
2	BMA	0	3	2	11,11,12	0.25	0	15,15,17	0.56	0
2	NAG	1	1	1,2	14,14,15	0.35	0	17,19,21	0.41	0
2	NAG	1	2	2	14,14,15	0.19	0	17,19,21	0.51	0
2	BMA	1	3	2	11,11,12	0.59	0	15,15,17	0.69	0
6	NAG	2	1	6,1	14,14,15	0.38	0	17,19,21	0.55	0
6	NAG	2	2	6	14,14,15	0.38	0	17,19,21	0.79	1 (5%)
6	BMA	2	3	6	11,11,12	0.29	0	15,15,17	0.59	0
6	MAN	2	4	6	11,11,12	0.25	0	15,15,17	0.58	0
6	MAN	2	5	6	11,11,12	0.30	0	15,15,17	0.68	0
6	MAN	2	6	6	11,11,12	0.31	0	15,15,17	0.82	1 (6%)
5	NAG	3	1	1,5	14,14,15	0.59	0	17,19,21	0.71	0
5	NAG	3	2	5	14,14,15	0.38	0	17,19,21	1.25	2 (11%)
5	BMA	3	3	5	11,11,12	0.35	0	15,15,17	0.52	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MAN	3	4	5	11,11,12	0.31	0	15,15,17	0.53	0
2	NAG	4	1	1,2	14,14,15	0.37	0	17,19,21	0.62	0
2	NAG	4	2	2	14,14,15	0.38	0	17,19,21	0.82	1 (5%)
2	BMA	4	3	2	11,11,12	0.26	0	15,15,17	0.56	0
4	NAG	5	1	1,4	14,14,15	0.29	0	17,19,21	0.52	0
4	NAG	5	2	4	14,14,15	0.36	0	17,19,21	0.75	1 (5%)
4	BMA	5	3	4	11,11,12	0.30	0	15,15,17	0.52	0
4	MAN	5	4	4	11,11,12	0.27	0	15,15,17	0.54	0
7	NAG	6	1	7,1	14,14,15	0.37	0	17,19,21	0.44	0
7	NAG	6	2	7	14,14,15	0.38	0	17,19,21	0.84	1 (5%)
7	BMA	6	3	7	11,11,12	0.27	0	15,15,17	0.77	1 (6%)
7	MAN	6	4	7	11,11,12	0.29	0	15,15,17	0.53	0
7	MAN	6	5	7	11,11,12	0.30	0	15,15,17	0.56	0
11	NAG	7	1	1,11	14,14,15	0.67	1 (7%)	17,19,21	0.79	0
11	NAG	7	2	11	14,14,15	0.20	0	17,19,21	0.42	0
11	BMA	7	3	11	11,11,12	0.91	1 (9%)	15,15,17	1.09	2 (13%)
11	MAN	7	4	11	11,11,12	1.33	2 (18%)	15,15,17	1.85	2 (13%)
11	MAN	7	5	11	11,11,12	0.83	0	15,15,17	1.05	2 (13%)
11	MAN	7	6	11	11,11,12	0.59	0	15,15,17	1.22	2 (13%)
11	MAN	7	7	11	11,11,12	0.70	0	15,15,17	1.08	2 (13%)
2	NAG	8	1	1,2	14,14,15	0.39	0	17,19,21	0.85	2 (11%)
2	NAG	8	2	2	14,14,15	0.52	0	17,19,21	0.80	1 (5%)
2	BMA	8	3	2	11,11,12	0.59	0	15,15,17	1.23	2 (13%)
5	NAG	9	1	5	14,14,15	0.26	0	17,19,21	1.06	2 (11%)
5	NAG	9	2	5	14,14,15	0.33	0	17,19,21	1.58	4 (23%)
5	BMA	9	3	5	11,11,12	0.20	0	15,15,17	0.57	0
5	MAN	9	4	5	11,11,12	0.22	0	15,15,17	0.54	0
2	NAG	AA	1	1,2	14,14,15	0.38	0	17,19,21	0.47	0
2	NAG	AA	2	2	14,14,15	0.38	0	17,19,21	0.60	0
2	BMA	AA	3	2	11,11,12	0.27	0	15,15,17	0.49	0
2	NAG	BA	1	1,2	14,14,15	0.31	0	17,19,21	0.79	0
2	NAG	BA	2	2	14,14,15	0.30	0	17,19,21	0.52	0
2	BMA	BA	3	2	11,11,12	0.21	0	15,15,17	0.55	0
9	NAG	CA	1	1,9	14,14,15	1.13	1 (7%)	17,19,21	2.23	5 (29%)
9	NAG	CA	2	9	14,14,15	0.30	0	17,19,21	0.58	0
9	BMA	CA	3	9	11,11,12	0.21	0	15,15,17	0.63	0
9	FUC	CA	4	9	10,10,11	0.27	0	14,14,16	0.53	0
12	NAG	DA	1	1,12	14,14,15	0.34	0	17,19,21	0.52	0
12	NAG	DA	2	12	14,14,15	0.26	0	17,19,21	0.69	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
12	FUC	DA	3	12	10,10,11	0.27	0	14,14,16	0.53	0
3	NAG	EA	1	1,3	14,14,15	0.30	0	17,19,21	0.61	0
3	NAG	EA	2	3	14,14,15	0.40	0	17,19,21	2.00	3 (17%)
12	NAG	FA	1	1,12	14,14,15	0.30	0	17,19,21	1.19	1 (5%)
12	NAG	FA	2	12	14,14,15	0.30	0	17,19,21	0.58	0
12	FUC	FA	3	12	10,10,11	0.23	0	14,14,16	0.52	0
2	NAG	G	1	1,2	14,14,15	0.33	0	17,19,21	0.56	0
2	NAG	G	2	2	14,14,15	0.30	0	17,19,21	0.60	0
2	BMA	G	3	2	11,11,12	0.25	0	15,15,17	0.56	0
3	NAG	H	1	1,3	14,14,15	0.34	0	17,19,21	0.92	1 (5%)
3	NAG	H	2	3	14,14,15	0.42	0	17,19,21	2.01	4 (23%)
2	NAG	I	1	1,2	14,14,15	0.30	0	17,19,21	1.31	2 (11%)
2	NAG	I	2	2	14,14,15	0.30	0	17,19,21	0.63	0
2	BMA	I	3	2	11,11,12	0.57	0	15,15,17	2.37	2 (13%)
3	NAG	J	1	1,3	14,14,15	0.30	0	17,19,21	0.53	0
3	NAG	J	2	3	14,14,15	0.32	0	17,19,21	0.65	0
4	NAG	K	1	1,4	14,14,15	0.39	0	17,19,21	0.53	0
4	NAG	K	2	4	14,14,15	0.39	0	17,19,21	1.19	2 (11%)
4	BMA	K	3	4	11,11,12	0.23	0	15,15,17	0.58	0
4	MAN	K	4	4	11,11,12	0.31	0	15,15,17	0.54	0
5	NAG	L	1	1,5	14,14,15	0.38	0	17,19,21	0.96	2 (11%)
5	NAG	L	2	5	14,14,15	0.38	0	17,19,21	1.33	2 (11%)
5	BMA	L	3	5	11,11,12	0.28	0	15,15,17	0.55	0
5	MAN	L	4	5	11,11,12	0.30	0	15,15,17	0.55	0
3	NAG	M	1	1,3	14,14,15	0.31	0	17,19,21	0.73	0
3	NAG	M	2	3	14,14,15	0.29	0	17,19,21	0.52	0
2	NAG	N	1	1,2	14,14,15	0.47	0	17,19,21	1.48	4 (23%)
2	NAG	N	2	2	14,14,15	0.38	0	17,19,21	1.00	1 (5%)
2	BMA	N	3	2	11,11,12	0.33	0	15,15,17	1.78	3 (20%)
3	NAG	O	1	1,3	14,14,15	0.50	0	17,19,21	0.71	0
3	NAG	O	2	3	14,14,15	0.20	0	17,19,21	0.84	0
3	NAG	P	1	1,3	14,14,15	0.32	0	17,19,21	0.95	1 (5%)
3	NAG	P	2	3	14,14,15	0.27	0	17,19,21	0.56	0
6	NAG	Q	1	6,1	14,14,15	1.28	1 (7%)	17,19,21	3.33	6 (35%)
6	NAG	Q	2	6	14,14,15	0.28	0	17,19,21	1.18	2 (11%)
6	BMA	Q	3	6	11,11,12	0.23	0	15,15,17	0.71	0
6	MAN	Q	4	6	11,11,12	0.16	0	15,15,17	0.64	0
6	MAN	Q	5	6	11,11,12	0.21	0	15,15,17	0.55	0
6	MAN	Q	6	6	11,11,12	0.20	0	15,15,17	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	R	1	1,2	14,14,15	0.58	0	17,19,21	0.70	0
2	NAG	R	2	2	14,14,15	0.25	0	17,19,21	0.85	0
2	BMA	R	3	2	11,11,12	1.26	1 (9%)	15,15,17	1.76	3 (20%)
3	NAG	S	1	1,3	14,14,15	0.30	0	17,19,21	0.52	0
3	NAG	S	2	3	14,14,15	0.24	0	17,19,21	0.55	0
2	NAG	T	1	1,2	14,14,15	0.33	0	17,19,21	0.69	0
2	NAG	T	2	2	14,14,15	0.24	0	17,19,21	0.89	1 (5%)
2	BMA	T	3	2	11,11,12	0.22	0	15,15,17	0.57	0
5	NAG	U	1	1,5	14,14,15	0.35	0	17,19,21	0.50	0
5	NAG	U	2	5	14,14,15	0.21	0	17,19,21	0.74	0
5	BMA	U	3	5	11,11,12	0.89	1 (9%)	15,15,17	0.88	0
5	MAN	U	4	5	11,11,12	0.70	1 (9%)	15,15,17	1.35	2 (13%)
7	NAG	V	1	7,1	14,14,15	0.40	0	17,19,21	0.88	1 (5%)
7	NAG	V	2	7	14,14,15	0.38	0	17,19,21	0.59	0
7	BMA	V	3	7	11,11,12	0.26	0	15,15,17	0.60	0
7	MAN	V	4	7	11,11,12	0.30	0	15,15,17	0.67	0
7	MAN	V	5	7	11,11,12	0.28	0	15,15,17	0.59	0
8	NAG	W	1	1,8	14,14,15	0.38	0	17,19,21	0.96	2 (11%)
8	NAG	W	2	8	14,14,15	0.38	0	17,19,21	0.48	0
8	MAN	W	3	8	11,11,12	0.30	0	15,15,17	0.52	0
8	BMA	W	4	8	11,11,12	0.26	0	15,15,17	0.58	0
5	NAG	X	1	1,5	14,14,15	0.38	0	17,19,21	0.85	1 (5%)
5	NAG	X	2	5	14,14,15	0.38	0	17,19,21	0.85	1 (5%)
5	BMA	X	3	5	11,11,12	0.27	0	15,15,17	0.58	0
5	MAN	X	4	5	11,11,12	0.31	0	15,15,17	0.56	0
2	NAG	Y	1	1,2	14,14,15	0.37	0	17,19,21	0.51	0
2	NAG	Y	2	2	14,14,15	0.37	0	17,19,21	0.79	1 (5%)
2	BMA	Y	3	2	11,11,12	0.27	0	15,15,17	0.54	0
2	NAG	Z	1	1,2	14,14,15	0.29	0	17,19,21	0.46	0
2	NAG	Z	2	2	14,14,15	0.27	0	17,19,21	0.49	0
2	BMA	Z	3	2	11,11,12	0.94	0	15,15,17	1.10	2 (13%)
9	NAG	a	1	1,9	14,14,15	1.13	1 (7%)	17,19,21	2.23	5 (29%)
9	NAG	a	2	9	14,14,15	0.34	0	17,19,21	0.58	0
9	BMA	a	3	9	11,11,12	0.23	0	15,15,17	0.63	0
9	FUC	a	4	9	10,10,11	0.25	0	14,14,16	0.52	0
3	NAG	b	1	1,3	14,14,15	0.35	0	17,19,21	0.91	1 (5%)
3	NAG	b	2	3	14,14,15	0.43	0	17,19,21	2.00	4 (23%)
2	NAG	c	1	1,2	14,14,15	0.29	0	17,19,21	0.46	0
2	NAG	c	2	2	14,14,15	0.28	0	17,19,21	0.57	0
2	BMA	c	3	2	11,11,12	0.24	0	15,15,17	0.60	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	d	1	1,2	14,14,15	0.26	0	17,19,21	0.49	0
2	NAG	d	2	2	14,14,15	0.26	0	17,19,21	0.86	1 (5%)
2	BMA	d	3	2	11,11,12	0.23	0	15,15,17	0.53	0
7	NAG	e	1	7,1	14,14,15	0.37	0	17,19,21	0.76	1 (5%)
7	NAG	e	2	7	14,14,15	0.35	0	17,19,21	1.03	1 (5%)
7	BMA	e	3	7	11,11,12	0.28	0	15,15,17	0.65	0
7	MAN	e	4	7	11,11,12	0.42	0	15,15,17	1.25	2 (13%)
7	MAN	e	5	7	11,11,12	0.30	0	15,15,17	0.53	0
7	NAG	f	1	7,1	14,14,15	0.27	0	17,19,21	0.86	1 (5%)
7	NAG	f	2	7	14,14,15	0.30	0	17,19,21	0.63	0
7	BMA	f	3	7	11,11,12	0.24	0	15,15,17	0.64	0
7	MAN	f	4	7	11,11,12	0.20	0	15,15,17	0.56	0
7	MAN	f	5	7	11,11,12	0.23	0	15,15,17	0.54	0
2	NAG	g	1	1,2	14,14,15	0.39	0	17,19,21	0.72	0
2	NAG	g	2	2	14,14,15	0.39	0	17,19,21	0.95	2 (11%)
2	BMA	g	3	2	11,11,12	0.27	0	15,15,17	0.56	0
3	NAG	h	1	1,3	14,14,15	0.35	0	17,19,21	1.61	2 (11%)
3	NAG	h	2	3	14,14,15	0.30	0	17,19,21	0.65	0
4	NAG	i	1	1,4	14,14,15	0.36	0	17,19,21	0.55	0
4	NAG	i	2	4	14,14,15	0.41	0	17,19,21	0.54	0
4	BMA	i	3	4	11,11,12	0.26	0	15,15,17	0.54	0
4	MAN	i	4	4	11,11,12	0.30	0	15,15,17	0.54	0
2	NAG	j	1	1,2	14,14,15	0.51	0	17,19,21	1.70	5 (29%)
2	NAG	j	2	2	14,14,15	0.28	0	17,19,21	0.86	2 (11%)
2	BMA	j	3	2	11,11,12	0.31	0	15,15,17	0.98	0
2	NAG	k	1	1,2	14,14,15	0.28	0	17,19,21	0.84	0
2	NAG	k	2	2	14,14,15	0.31	0	17,19,21	0.65	0
2	BMA	k	3	2	11,11,12	0.20	0	15,15,17	0.60	0
10	NAG	l	1	1,10	14,14,15	0.39	0	17,19,21	0.59	0
10	NAG	l	2	10	14,14,15	0.39	0	17,19,21	0.50	0
10	BMA	l	3	10	11,11,12	0.34	0	15,15,17	0.52	0
10	MAN	l	4	10	11,11,12	0.31	0	15,15,17	0.51	0
10	MAN	l	5	10	11,11,12	0.28	0	15,15,17	0.58	0
10	MAN	l	6	10	11,11,12	0.29	0	15,15,17	0.72	0
10	MAN	l	7	10	11,11,12	0.29	0	15,15,17	0.53	0
10	MAN	l	8	10	11,11,12	0.24	0	15,15,17	0.53	0
5	NAG	m	1	1,5	14,14,15	0.38	0	17,19,21	0.73	1 (5%)
5	NAG	m	2	5	14,14,15	0.38	0	17,19,21	0.77	1 (5%)
5	BMA	m	3	5	11,11,12	0.28	0	15,15,17	0.51	0
5	MAN	m	4	5	11,11,12	0.28	0	15,15,17	0.70	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	n	1	1,2	14,14,15	0.31	0	17,19,21	0.58	0
2	NAG	n	2	2	14,14,15	0.24	0	17,19,21	0.76	0
2	BMA	n	3	2	11,11,12	0.22	0	15,15,17	0.58	0
2	NAG	o	1	1,2	14,14,15	0.29	0	17,19,21	0.62	0
2	NAG	o	2	2	14,14,15	0.31	0	17,19,21	0.56	0
2	BMA	o	3	2	11,11,12	0.22	0	15,15,17	0.56	0
4	NAG	p	1	1,4	14,14,15	0.28	0	17,19,21	0.41	0
4	NAG	p	2	4	14,14,15	0.32	0	17,19,21	0.68	0
4	BMA	p	3	4	11,11,12	0.24	0	15,15,17	0.63	0
4	MAN	p	4	4	11,11,12	0.20	0	15,15,17	0.67	0
11	NAG	q	1	1,11	14,14,15	0.71	1 (7%)	17,19,21	0.78	0
11	NAG	q	2	11	14,14,15	0.20	0	17,19,21	0.40	0
11	BMA	q	3	11	11,11,12	0.93	1 (9%)	15,15,17	1.07	1 (6%)
11	MAN	q	4	11	11,11,12	1.36	2 (18%)	15,15,17	1.84	2 (13%)
11	MAN	q	5	11	11,11,12	0.84	1 (9%)	15,15,17	1.04	2 (13%)
11	MAN	q	6	11	11,11,12	0.60	0	15,15,17	1.23	2 (13%)
11	MAN	q	7	11	11,11,12	0.71	0	15,15,17	1.08	2 (13%)
5	NAG	r	1	1,5	14,14,15	0.35	0	17,19,21	0.85	2 (11%)
5	NAG	r	2	5	14,14,15	0.49	0	17,19,21	0.81	1 (5%)
5	BMA	r	3	5	11,11,12	1.03	1 (9%)	15,15,17	0.93	1 (6%)
5	MAN	r	4	5	11,11,12	0.76	0	15,15,17	1.26	2 (13%)
2	NAG	s	1	1,2	14,14,15	0.28	0	17,19,21	0.97	2 (11%)
2	NAG	s	2	2	14,14,15	0.32	0	17,19,21	0.61	0
2	BMA	s	3	2	11,11,12	0.22	0	15,15,17	0.56	0
2	NAG	t	1	1,2	14,14,15	0.29	0	17,19,21	0.62	0
2	NAG	t	2	2	14,14,15	0.26	0	17,19,21	1.01	1 (5%)
2	BMA	t	3	2	11,11,12	0.25	0	15,15,17	0.56	0
2	NAG	u	1	1,2	14,14,15	0.32	0	17,19,21	0.80	1 (5%)
2	NAG	u	2	2	14,14,15	0.24	0	17,19,21	0.85	1 (5%)
2	BMA	u	3	2	11,11,12	0.23	0	15,15,17	0.58	0
3	NAG	v	1	1,3	14,14,15	0.49	0	17,19,21	0.51	0
3	NAG	v	2	3	14,14,15	0.32	0	17,19,21	0.42	0
7	NAG	w	1	7,1	14,14,15	0.37	0	17,19,21	0.75	1 (5%)
7	NAG	w	2	7	14,14,15	0.37	0	17,19,21	1.20	2 (11%)
7	BMA	w	3	7	11,11,12	0.29	0	15,15,17	0.66	0
7	MAN	w	4	7	11,11,12	0.29	0	15,15,17	0.55	0
7	MAN	w	5	7	11,11,12	0.27	0	15,15,17	0.57	0
2	NAG	x	1	1,2	14,14,15	0.44	0	17,19,21	0.54	0
2	NAG	x	2	2	14,14,15	0.36	0	17,19,21	0.80	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BMA	x	3	2	11,11,12	1.09	0	15,15,17	1.13	2 (13%)
2	NAG	y	1	1,2	14,14,15	0.40	0	17,19,21	0.55	0
2	NAG	y	2	2	14,14,15	0.40	0	17,19,21	1.52	2 (11%)
2	BMA	y	3	2	11,11,12	0.27	0	15,15,17	0.58	0
2	NAG	z	1	1,2	14,14,15	0.29	0	17,19,21	0.58	0
2	NAG	z	2	2	14,14,15	0.33	0	17,19,21	1.60	4 (23%)
2	BMA	z	3	2	11,11,12	0.22	0	15,15,17	0.54	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	0	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	0	2	2	-	4/6/23/26	0/1/1/1
2	BMA	0	3	2	-	0/2/19/22	0/1/1/1
2	NAG	1	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	1	2	2	-	1/6/23/26	0/1/1/1
2	BMA	1	3	2	-	0/2/19/22	0/1/1/1
6	NAG	2	1	6,1	-	0/6/23/26	0/1/1/1
6	NAG	2	2	6	-	2/6/23/26	0/1/1/1
6	BMA	2	3	6	-	2/2/19/22	0/1/1/1
6	MAN	2	4	6	-	1/2/19/22	0/1/1/1
6	MAN	2	5	6	-	0/2/19/22	0/1/1/1
6	MAN	2	6	6	-	0/2/19/22	0/1/1/1
5	NAG	3	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	3	2	5	-	3/6/23/26	0/1/1/1
5	BMA	3	3	5	-	0/2/19/22	0/1/1/1
5	MAN	3	4	5	-	0/2/19/22	0/1/1/1
2	NAG	4	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	4	2	2	-	1/6/23/26	0/1/1/1
2	BMA	4	3	2	-	0/2/19/22	0/1/1/1
4	NAG	5	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	5	2	4	-	3/6/23/26	0/1/1/1
4	BMA	5	3	4	-	0/2/19/22	0/1/1/1
4	MAN	5	4	4	-	1/2/19/22	0/1/1/1
7	NAG	6	1	7,1	-	0/6/23/26	0/1/1/1
7	NAG	6	2	7	-	1/6/23/26	0/1/1/1
7	BMA	6	3	7	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	MAN	6	4	7	-	0/2/19/22	0/1/1/1
7	MAN	6	5	7	-	0/2/19/22	0/1/1/1
11	NAG	7	1	1,11	-	0/6/23/26	0/1/1/1
11	NAG	7	2	11	-	0/6/23/26	0/1/1/1
11	BMA	7	3	11	-	2/2/19/22	0/1/1/1
11	MAN	7	4	11	-	0/2/19/22	0/1/1/1
11	MAN	7	5	11	-	0/2/19/22	0/1/1/1
11	MAN	7	6	11	-	0/2/19/22	0/1/1/1
11	MAN	7	7	11	-	0/2/19/22	0/1/1/1
2	NAG	8	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	8	2	2	-	2/6/23/26	0/1/1/1
2	BMA	8	3	2	-	0/2/19/22	0/1/1/1
5	NAG	9	1	5	-	3/6/23/26	0/1/1/1
5	NAG	9	2	5	-	3/6/23/26	0/1/1/1
5	BMA	9	3	5	-	0/2/19/22	0/1/1/1
5	MAN	9	4	5	-	1/2/19/22	0/1/1/1
2	NAG	AA	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	AA	2	2	-	0/6/23/26	0/1/1/1
2	BMA	AA	3	2	-	1/2/19/22	0/1/1/1
2	NAG	BA	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	BA	2	2	-	0/6/23/26	0/1/1/1
2	BMA	BA	3	2	-	0/2/19/22	0/1/1/1
9	NAG	CA	1	1,9	-	0/6/23/26	0/1/1/1
9	NAG	CA	2	9	-	0/6/23/26	0/1/1/1
9	BMA	CA	3	9	-	0/2/19/22	0/1/1/1
9	FUC	CA	4	9	-	-	0/1/1/1
12	NAG	DA	1	1,12	-	0/6/23/26	0/1/1/1
12	NAG	DA	2	12	-	0/6/23/26	0/1/1/1
12	FUC	DA	3	12	-	-	0/1/1/1
3	NAG	EA	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	EA	2	3	-	1/6/23/26	0/1/1/1
12	NAG	FA	1	1,12	-	3/6/23/26	0/1/1/1
12	NAG	FA	2	12	-	2/6/23/26	0/1/1/1
12	FUC	FA	3	12	-	-	0/1/1/1
2	NAG	G	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	BMA	G	3	2	-	0/2/19/22	0/1/1/1
3	NAG	H	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	H	2	3	-	1/6/23/26	0/1/1/1
2	NAG	I	1	1,2	-	3/6/23/26	0/1/1/1

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	BMA	U	3	5	-	0/2/19/22	0/1/1/1
5	MAN	U	4	5	-	2/2/19/22	0/1/1/1
7	NAG	V	1	7,1	-	0/6/23/26	0/1/1/1
7	NAG	V	2	7	-	0/6/23/26	0/1/1/1
7	BMA	V	3	7	-	0/2/19/22	0/1/1/1
7	MAN	V	4	7	-	0/2/19/22	0/1/1/1
7	MAN	V	5	7	-	1/2/19/22	0/1/1/1
8	NAG	W	1	1,8	-	4/6/23/26	0/1/1/1
8	NAG	W	2	8	-	0/6/23/26	0/1/1/1
8	MAN	W	3	8	-	0/2/19/22	0/1/1/1
8	BMA	W	4	8	-	0/2/19/22	0/1/1/1
5	NAG	X	1	1,5	-	1/6/23/26	0/1/1/1
5	NAG	X	2	5	-	1/6/23/26	0/1/1/1
5	BMA	X	3	5	-	0/2/19/22	0/1/1/1
5	MAN	X	4	5	-	0/2/19/22	0/1/1/1
2	NAG	Y	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	Y	2	2	-	1/6/23/26	0/1/1/1
2	BMA	Y	3	2	-	0/2/19/22	0/1/1/1
2	NAG	Z	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	Z	2	2	-	2/6/23/26	0/1/1/1
2	BMA	Z	3	2	-	2/2/19/22	0/1/1/1
9	NAG	a	1	1,9	-	0/6/23/26	0/1/1/1
9	NAG	a	2	9	-	0/6/23/26	0/1/1/1
9	BMA	a	3	9	-	0/2/19/22	0/1/1/1
9	FUC	a	4	9	-	-	0/1/1/1
3	NAG	b	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	b	2	3	-	1/6/23/26	0/1/1/1
2	NAG	c	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	c	2	2	-	1/6/23/26	0/1/1/1
2	BMA	c	3	2	-	1/2/19/22	0/1/1/1
2	NAG	d	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	d	2	2	-	3/6/23/26	0/1/1/1
2	BMA	d	3	2	-	0/2/19/22	0/1/1/1
7	NAG	e	1	7,1	-	2/6/23/26	0/1/1/1
7	NAG	e	2	7	-	4/6/23/26	0/1/1/1
7	BMA	e	3	7	-	2/2/19/22	0/1/1/1
7	MAN	e	4	7	-	0/2/19/22	0/1/1/1
7	MAN	e	5	7	-	0/2/19/22	0/1/1/1
7	NAG	f	1	7,1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	f	2	7	-	2/6/23/26	0/1/1/1
7	BMA	f	3	7	-	0/2/19/22	0/1/1/1
7	MAN	f	4	7	-	1/2/19/22	0/1/1/1
7	MAN	f	5	7	-	0/2/19/22	0/1/1/1
2	NAG	g	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	g	2	2	-	5/6/23/26	0/1/1/1
2	BMA	g	3	2	-	0/2/19/22	0/1/1/1
3	NAG	h	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	h	2	3	-	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	-	1/6/23/26	0/1/1/1
4	BMA	i	3	4	-	0/2/19/22	0/1/1/1
4	MAN	i	4	4	-	0/2/19/22	0/1/1/1
2	NAG	j	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	j	2	2	-	3/6/23/26	0/1/1/1
2	BMA	j	3	2	-	0/2/19/22	0/1/1/1
2	NAG	k	1	1,2	-	5/6/23/26	0/1/1/1
2	NAG	k	2	2	-	2/6/23/26	0/1/1/1
2	BMA	k	3	2	-	0/2/19/22	0/1/1/1
10	NAG	l	1	1,10	-	0/6/23/26	0/1/1/1
10	NAG	l	2	10	-	0/6/23/26	0/1/1/1
10	BMA	l	3	10	-	0/2/19/22	0/1/1/1
10	MAN	l	4	10	-	0/2/19/22	0/1/1/1
10	MAN	l	5	10	-	0/2/19/22	0/1/1/1
10	MAN	l	6	10	-	1/2/19/22	0/1/1/1
10	MAN	l	7	10	-	0/2/19/22	0/1/1/1
10	MAN	l	8	10	-	0/2/19/22	0/1/1/1
5	NAG	m	1	1,5	-	1/6/23/26	0/1/1/1
5	NAG	m	2	5	-	1/6/23/26	0/1/1/1
5	BMA	m	3	5	-	0/2/19/22	0/1/1/1
5	MAN	m	4	5	-	0/2/19/22	0/1/1/1
2	NAG	n	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	n	2	2	-	3/6/23/26	0/1/1/1
2	BMA	n	3	2	-	0/2/19/22	0/1/1/1
2	NAG	o	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	o	2	2	-	0/6/23/26	0/1/1/1
2	BMA	o	3	2	-	0/2/19/22	0/1/1/1
4	NAG	p	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	p	2	4	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	BMA	p	3	4	-	2/2/19/22	0/1/1/1
4	MAN	p	4	4	-	0/2/19/22	0/1/1/1
11	NAG	q	1	1,11	-	0/6/23/26	0/1/1/1
11	NAG	q	2	11	-	0/6/23/26	0/1/1/1
11	BMA	q	3	11	-	2/2/19/22	0/1/1/1
11	MAN	q	4	11	-	0/2/19/22	0/1/1/1
11	MAN	q	5	11	-	0/2/19/22	0/1/1/1
11	MAN	q	6	11	-	0/2/19/22	0/1/1/1
11	MAN	q	7	11	-	0/2/19/22	0/1/1/1
5	NAG	r	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	r	2	5	-	2/6/23/26	0/1/1/1
5	BMA	r	3	5	-	0/2/19/22	0/1/1/1
5	MAN	r	4	5	-	0/2/19/22	0/1/1/1
2	NAG	s	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	s	2	2	-	0/6/23/26	0/1/1/1
2	BMA	s	3	2	-	0/2/19/22	0/1/1/1
2	NAG	t	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	t	2	2	-	3/6/23/26	0/1/1/1
2	BMA	t	3	2	-	2/2/19/22	0/1/1/1
2	NAG	u	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	u	2	2	-	4/6/23/26	0/1/1/1
2	BMA	u	3	2	-	0/2/19/22	0/1/1/1
3	NAG	v	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	v	2	3	-	0/6/23/26	0/1/1/1
7	NAG	w	1	7,1	-	1/6/23/26	0/1/1/1
7	NAG	w	2	7	-	3/6/23/26	0/1/1/1
7	BMA	w	3	7	-	0/2/19/22	0/1/1/1
7	MAN	w	4	7	-	1/2/19/22	0/1/1/1
7	MAN	w	5	7	-	0/2/19/22	0/1/1/1
2	NAG	x	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	x	2	2	-	2/6/23/26	0/1/1/1
2	BMA	x	3	2	-	1/2/19/22	0/1/1/1
2	NAG	y	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	y	2	2	-	3/6/23/26	0/1/1/1
2	BMA	y	3	2	-	0/2/19/22	0/1/1/1
2	NAG	z	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	z	2	2	-	3/6/23/26	0/1/1/1
2	BMA	z	3	2	-	0/2/19/22	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	Q	1	NAG	C1-C2	4.21	1.58	1.52
9	CA	1	NAG	C1-C2	4.01	1.57	1.52
9	a	1	NAG	C1-C2	4.00	1.57	1.52
2	R	3	BMA	O5-C5	3.42	1.50	1.43
11	7	4	MAN	C1-C2	3.23	1.59	1.52
11	q	4	MAN	C1-C2	3.20	1.59	1.52
5	r	3	BMA	C1-C2	2.89	1.59	1.52
11	q	4	MAN	O5-C1	2.53	1.47	1.43
11	q	1	NAG	O5-C1	-2.52	1.39	1.43
11	7	4	MAN	O5-C1	2.44	1.47	1.43
11	7	1	NAG	O5-C1	-2.35	1.39	1.43
5	U	3	BMA	O5-C1	-2.21	1.40	1.43
11	q	3	BMA	C4-C3	2.15	1.57	1.52
11	7	3	BMA	C4-C3	2.10	1.57	1.52
11	q	5	MAN	C1-C2	2.07	1.57	1.52
5	U	4	MAN	C1-C2	2.02	1.57	1.52

All (146) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	Q	1	NAG	C1-O5-C5	9.98	125.56	112.19
2	I	3	BMA	O2-C2-C3	8.27	127.27	110.15
9	a	1	NAG	O5-C1-C2	-7.47	99.74	111.29
9	CA	1	NAG	O5-C1-C2	-7.46	99.75	111.29
6	Q	1	NAG	O5-C1-C2	-6.64	101.02	111.29
3	H	2	NAG	O5-C5-C6	6.54	120.40	107.66
3	EA	2	NAG	O5-C5-C6	6.53	120.38	107.66
3	b	2	NAG	O5-C5-C6	6.53	120.38	107.66
11	7	4	MAN	C1-O5-C5	6.07	120.32	112.19
11	q	4	MAN	C1-O5-C5	6.00	120.23	112.19
2	N	3	BMA	C1-O5-C5	5.21	119.17	112.19
3	h	1	NAG	C2-N2-C7	4.63	129.10	122.90
2	R	3	BMA	C1-O5-C5	4.57	118.31	112.19
2	z	2	NAG	C1-O5-C5	4.51	118.23	112.19
5	9	2	NAG	C1-O5-C5	4.46	118.17	112.19
2	y	2	NAG	C2-N2-C7	4.12	128.42	122.90
5	U	4	MAN	C1-O5-C5	4.02	117.57	112.19
2	y	2	NAG	C1-C2-N2	4.00	116.73	110.43
5	r	4	MAN	C1-O5-C5	3.98	117.53	112.19
11	q	6	MAN	C1-O5-C5	3.82	117.31	112.19
2	I	1	NAG	C2-N2-C7	3.79	127.97	122.90
7	e	2	NAG	C2-N2-C7	3.78	127.96	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	7	6	MAN	C1-O5-C5	3.74	117.19	112.19
2	j	1	NAG	C4-C3-C2	-3.65	105.67	111.02
5	L	2	NAG	C1-C2-N2	3.47	115.91	110.43
6	Q	1	NAG	C2-N2-C7	-3.47	118.25	122.90
2	R	3	BMA	O3-C3-C2	3.45	117.09	110.05
5	3	2	NAG	C1-C2-N2	3.37	115.75	110.43
7	w	2	NAG	C1-C2-N2	3.37	115.75	110.43
3	h	1	NAG	C1-C2-N2	3.30	115.63	110.43
5	3	2	NAG	C2-N2-C7	3.27	127.28	122.90
2	N	1	NAG	O5-C1-C2	3.23	116.30	111.29
5	L	2	NAG	C2-N2-C7	3.21	127.20	122.90
9	a	1	NAG	C3-C4-C5	-3.21	104.42	110.23
4	K	2	NAG	C1-C2-N2	3.19	115.46	110.43
2	N	1	NAG	O4-C4-C5	3.17	117.14	109.32
2	j	1	NAG	O4-C4-C5	3.16	117.12	109.32
9	CA	1	NAG	C3-C4-C5	-3.13	104.55	110.23
2	N	2	NAG	C1-O5-C5	3.08	116.32	112.19
6	Q	2	NAG	C2-N2-C7	3.08	127.03	122.90
12	FA	1	NAG	C2-N2-C7	3.01	126.93	122.90
5	L	1	NAG	C1-C2-N2	2.98	115.14	110.43
2	N	3	BMA	C1-C2-C3	2.95	113.95	109.64
11	q	7	MAN	C1-O5-C5	2.93	116.12	112.19
11	7	7	MAN	C1-O5-C5	2.92	116.10	112.19
7	e	4	MAN	C1-O5-C5	2.90	116.08	112.19
7	e	4	MAN	O2-C2-C3	2.88	116.12	110.15
2	N	1	NAG	C4-C3-C2	-2.84	106.86	111.02
6	Q	2	NAG	C1-C2-N2	-2.84	105.96	110.43
11	q	5	MAN	C1-O5-C5	2.82	115.96	112.19
2	j	1	NAG	O5-C1-C2	2.80	115.62	111.29
4	K	2	NAG	C2-N2-C7	2.79	126.64	122.90
3	P	1	NAG	C2-N2-C7	2.79	126.63	122.90
11	7	5	MAN	C1-O5-C5	2.76	115.89	112.19
3	H	2	NAG	C4-C3-C2	2.75	115.05	111.02
2	x	3	BMA	O3-C3-C4	2.74	116.83	110.38
3	b	2	NAG	C4-C3-C2	2.73	115.03	111.02
3	EA	2	NAG	C4-C3-C2	2.71	114.98	111.02
7	f	1	NAG	C1-O5-C5	2.69	115.80	112.19
2	z	2	NAG	O5-C1-C2	2.68	115.44	111.29
6	Q	1	NAG	O4-C4-C3	-2.68	104.07	110.38
6	Q	1	NAG	O5-C5-C6	-2.67	102.47	107.66
5	X	1	NAG	C2-N2-C7	2.66	126.47	122.90
7	w	2	NAG	C2-N2-C7	2.66	126.47	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	1	NAG	C1-C2-N2	2.66	114.62	110.43
8	W	1	NAG	C1-C2-N2	2.66	114.62	110.43
7	6	2	NAG	C2-N2-C7	2.64	126.43	122.90
2	8	3	BMA	O2-C2-C3	-2.63	104.70	110.15
2	j	1	NAG	O5-C5-C4	-2.59	104.53	110.83
11	q	4	MAN	C1-C2-C3	2.56	113.38	109.64
9	a	1	NAG	C1-O5-C5	2.56	115.62	112.19
3	b	1	NAG	C1-O5-C5	2.54	115.59	112.19
9	CA	1	NAG	C1-O5-C5	2.54	115.59	112.19
3	H	1	NAG	C1-O5-C5	2.53	115.58	112.19
2	0	2	NAG	C1-C2-N2	2.53	114.42	110.43
5	9	1	NAG	C1-O5-C5	2.52	115.56	112.19
4	5	2	NAG	C2-N2-C7	2.52	126.27	122.90
11	7	4	MAN	C1-C2-C3	2.51	113.31	109.64
5	9	2	NAG	C3-C4-C5	2.50	114.76	110.23
6	Q	1	NAG	C4-C3-C2	2.47	114.64	111.02
2	Y	2	NAG	C2-N2-C7	2.47	126.21	122.90
7	e	1	NAG	C2-N2-C7	2.47	126.21	122.90
5	9	2	NAG	O5-C1-C2	2.47	115.11	111.29
2	R	3	BMA	C3-C4-C5	-2.46	105.78	110.23
2	4	2	NAG	C2-N2-C7	2.46	126.19	122.90
2	d	2	NAG	C2-N2-C7	2.45	126.18	122.90
3	EA	2	NAG	O3-C3-C4	2.44	116.14	110.38
5	m	2	NAG	C2-N2-C7	2.42	126.14	122.90
2	N	1	NAG	O5-C5-C4	-2.42	104.95	110.83
3	H	2	NAG	O3-C3-C4	2.41	116.07	110.38
3	b	2	NAG	O3-C3-C4	2.41	116.06	110.38
2	N	3	BMA	C3-C4-C5	2.40	114.58	110.23
5	X	2	NAG	C2-N2-C7	2.39	126.11	122.90
5	m	1	NAG	C2-N2-C7	2.39	126.10	122.90
5	9	1	NAG	C2-N2-C7	2.37	126.08	122.90
2	z	2	NAG	C3-C4-C5	2.36	114.51	110.23
6	2	6	MAN	C1-O5-C5	2.35	115.34	112.19
2	I	3	BMA	C2-C3-C4	2.35	114.99	110.86
2	Z	3	BMA	O2-C2-C3	-2.34	105.31	110.15
2	8	3	BMA	C2-C3-C4	2.34	114.97	110.86
7	w	1	NAG	C2-N2-C7	2.33	126.03	122.90
2	s	1	NAG	C2-N2-C7	2.31	126.00	122.90
2	T	2	NAG	C1-O5-C5	2.31	115.28	112.19
5	r	3	BMA	O2-C2-C3	-2.30	105.39	110.15
2	g	2	NAG	C1-C2-N2	2.28	114.02	110.43
2	u	2	NAG	C1-O5-C5	2.26	115.22	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	g	2	NAG	C2-N2-C7	2.25	125.92	122.90
11	q	6	MAN	O2-C2-C3	-2.25	105.50	110.15
11	7	6	MAN	O2-C2-C3	-2.24	105.51	110.15
2	j	1	NAG	C3-C4-C5	2.23	114.28	110.23
6	2	2	NAG	C2-N2-C7	2.23	125.89	122.90
5	U	4	MAN	O2-C2-C3	-2.23	105.54	110.15
2	x	3	BMA	C1-O5-C5	2.23	115.17	112.19
2	t	2	NAG	C2-N2-C7	2.22	125.88	122.90
11	7	7	MAN	O2-C2-C3	-2.21	105.57	110.15
11	q	7	MAN	O2-C2-C3	-2.18	105.64	110.15
5	r	1	NAG	O4-C4-C5	-2.18	103.96	109.32
2	8	1	NAG	O4-C4-C5	-2.17	103.97	109.32
2	j	2	NAG	C1-O5-C5	2.17	115.10	112.19
7	6	3	BMA	C1-O5-C5	2.14	115.05	112.19
9	CA	1	NAG	C1-C2-N2	2.13	113.79	110.43
11	7	5	MAN	O2-C2-C3	-2.13	105.75	110.15
5	9	2	NAG	O5-C5-C6	2.12	111.79	107.66
11	q	3	BMA	C1-C2-C3	-2.12	106.56	109.64
2	u	1	NAG	C1-O5-C5	2.11	115.01	112.19
5	m	4	MAN	C1-O5-C5	2.10	115.00	112.19
5	L	1	NAG	C2-N2-C7	2.10	125.71	122.90
9	a	1	NAG	C1-C2-N2	2.09	113.73	110.43
11	q	5	MAN	O2-C2-C3	-2.09	105.83	110.15
2	8	2	NAG	C1-O5-C5	2.09	114.98	112.19
5	r	2	NAG	C1-O5-C5	2.08	114.98	112.19
2	Z	3	BMA	C1-C2-C3	-2.08	106.61	109.64
5	r	4	MAN	O2-C2-C3	-2.08	105.84	110.15
9	CA	1	NAG	C4-C3-C2	-2.08	107.97	111.02
11	7	3	BMA	C1-O5-C5	2.08	114.97	112.19
7	V	1	NAG	O4-C4-C3	-2.06	105.52	110.38
11	7	3	BMA	C1-C2-C3	-2.06	106.65	109.64
9	a	1	NAG	C4-C3-C2	-2.05	108.01	111.02
8	W	1	NAG	C2-N2-C7	2.03	125.62	122.90
2	s	1	NAG	C1-O5-C5	2.03	114.90	112.19
3	H	2	NAG	O4-C4-C5	2.03	114.32	109.32
2	8	1	NAG	O4-C4-C3	2.02	115.14	110.38
2	j	2	NAG	C2-N2-C7	2.02	125.60	122.90
3	b	2	NAG	O4-C4-C5	2.02	114.29	109.32
5	r	1	NAG	O4-C4-C3	2.01	115.12	110.38
2	z	2	NAG	O5-C5-C6	2.01	111.58	107.66

There are no chirality outliers.

All (241) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	I	1	NAG	C1-C2-N2-C7
2	I	1	NAG	C8-C7-N2-C2
2	I	1	NAG	O7-C7-N2-C2
2	N	1	NAG	C8-C7-N2-C2
2	N	1	NAG	O7-C7-N2-C2
2	T	2	NAG	C3-C2-N2-C7
2	T	2	NAG	O7-C7-N2-C2
2	c	1	NAG	C8-C7-N2-C2
2	c	1	NAG	O7-C7-N2-C2
2	d	2	NAG	C1-C2-N2-C7
2	d	2	NAG	C8-C7-N2-C2
2	d	2	NAG	O7-C7-N2-C2
2	g	1	NAG	C8-C7-N2-C2
2	g	1	NAG	O7-C7-N2-C2
2	g	2	NAG	C1-C2-N2-C7
2	j	2	NAG	C1-C2-N2-C7
2	j	2	NAG	C8-C7-N2-C2
2	j	2	NAG	O7-C7-N2-C2
2	k	1	NAG	C1-C2-N2-C7
2	k	1	NAG	C8-C7-N2-C2
2	k	1	NAG	O7-C7-N2-C2
2	n	2	NAG	C1-C2-N2-C7
2	n	2	NAG	C8-C7-N2-C2
2	n	2	NAG	O7-C7-N2-C2
2	t	2	NAG	C1-C2-N2-C7
2	t	2	NAG	O7-C7-N2-C2
2	u	1	NAG	C8-C7-N2-C2
2	u	1	NAG	O7-C7-N2-C2
2	u	2	NAG	C3-C2-N2-C7
2	u	2	NAG	O7-C7-N2-C2
2	y	2	NAG	C1-C2-N2-C7
2	y	2	NAG	C8-C7-N2-C2
2	y	2	NAG	O7-C7-N2-C2
2	z	2	NAG	C8-C7-N2-C2
2	z	2	NAG	O7-C7-N2-C2
2	0	2	NAG	C1-C2-N2-C7
2	0	2	NAG	C8-C7-N2-C2
2	0	2	NAG	O7-C7-N2-C2
3	J	1	NAG	O7-C7-N2-C2
3	J	2	NAG	C8-C7-N2-C2
3	J	2	NAG	O7-C7-N2-C2
3	P	1	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	P	1	NAG	O7-C7-N2-C2
3	h	1	NAG	C1-C2-N2-C7
4	K	2	NAG	C1-C2-N2-C7
5	L	1	NAG	C1-C2-N2-C7
5	L	1	NAG	C8-C7-N2-C2
5	L	1	NAG	O7-C7-N2-C2
5	L	2	NAG	C1-C2-N2-C7
5	3	2	NAG	C1-C2-N2-C7
5	9	1	NAG	C1-C2-N2-C7
6	Q	2	NAG	C3-C2-N2-C7
6	Q	2	NAG	C8-C7-N2-C2
6	Q	2	NAG	O7-C7-N2-C2
7	e	2	NAG	C3-C2-N2-C7
7	e	2	NAG	C8-C7-N2-C2
7	e	2	NAG	O7-C7-N2-C2
7	f	2	NAG	C8-C7-N2-C2
7	f	2	NAG	O7-C7-N2-C2
7	w	2	NAG	C1-C2-N2-C7
7	w	2	NAG	O7-C7-N2-C2
12	FA	1	NAG	C1-C2-N2-C7
12	FA	1	NAG	O7-C7-N2-C2
2	T	2	NAG	C8-C7-N2-C2
2	s	1	NAG	C8-C7-N2-C2
2	s	1	NAG	O7-C7-N2-C2
2	t	2	NAG	C8-C7-N2-C2
2	u	2	NAG	C8-C7-N2-C2
3	J	1	NAG	C8-C7-N2-C2
4	K	2	NAG	C8-C7-N2-C2
4	K	2	NAG	O7-C7-N2-C2
5	9	2	NAG	C8-C7-N2-C2
5	9	2	NAG	O7-C7-N2-C2
7	w	2	NAG	C8-C7-N2-C2
12	FA	1	NAG	C8-C7-N2-C2
5	U	4	MAN	O5-C5-C6-O6
2	x	1	NAG	C4-C5-C6-O6
3	h	1	NAG	C8-C7-N2-C2
3	h	1	NAG	O7-C7-N2-C2
5	L	2	NAG	C8-C7-N2-C2
5	L	2	NAG	O7-C7-N2-C2
5	3	2	NAG	O7-C7-N2-C2
5	9	1	NAG	O7-C7-N2-C2
4	p	3	BMA	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	6	3	BMA	O5-C5-C6-O6
11	q	3	BMA	O5-C5-C6-O6
11	7	3	BMA	O5-C5-C6-O6
2	x	2	NAG	O5-C5-C6-O6
2	Z	3	BMA	C4-C5-C6-O6
2	R	2	NAG	O5-C5-C6-O6
3	O	2	NAG	O5-C5-C6-O6
4	K	3	BMA	O5-C5-C6-O6
4	p	2	NAG	C8-C7-N2-C2
5	3	2	NAG	C8-C7-N2-C2
5	9	1	NAG	C8-C7-N2-C2
12	FA	2	NAG	C8-C7-N2-C2
2	Z	2	NAG	O5-C5-C6-O6
2	Z	3	BMA	O5-C5-C6-O6
2	t	3	BMA	O5-C5-C6-O6
6	2	3	BMA	O5-C5-C6-O6
2	t	3	BMA	C4-C5-C6-O6
3	O	2	NAG	C4-C5-C6-O6
2	x	1	NAG	O5-C5-C6-O6
2	R	2	NAG	C4-C5-C6-O6
5	U	4	MAN	C4-C5-C6-O6
11	q	3	BMA	C4-C5-C6-O6
4	p	3	BMA	C4-C5-C6-O6
11	7	3	BMA	C4-C5-C6-O6
2	T	1	NAG	C8-C7-N2-C2
2	z	1	NAG	C8-C7-N2-C2
3	H	1	NAG	C8-C7-N2-C2
3	b	1	NAG	C8-C7-N2-C2
7	f	1	NAG	C8-C7-N2-C2
12	FA	2	NAG	O7-C7-N2-C2
4	K	3	BMA	C4-C5-C6-O6
2	Z	2	NAG	C4-C5-C6-O6
2	8	2	NAG	C4-C5-C6-O6
5	r	2	NAG	C4-C5-C6-O6
2	Z	1	NAG	O5-C5-C6-O6
2	x	2	NAG	C4-C5-C6-O6
7	6	3	BMA	C4-C5-C6-O6
7	e	3	BMA	O5-C5-C6-O6
2	T	1	NAG	O7-C7-N2-C2
2	d	1	NAG	C8-C7-N2-C2
2	j	1	NAG	C8-C7-N2-C2
2	k	2	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	t	1	NAG	C8-C7-N2-C2
2	t	1	NAG	O7-C7-N2-C2
2	z	1	NAG	O7-C7-N2-C2
2	AA	1	NAG	C8-C7-N2-C2
2	AA	1	NAG	O7-C7-N2-C2
3	H	1	NAG	O7-C7-N2-C2
3	P	2	NAG	C8-C7-N2-C2
3	b	1	NAG	O7-C7-N2-C2
3	h	2	NAG	C8-C7-N2-C2
4	p	2	NAG	O7-C7-N2-C2
7	f	1	NAG	O7-C7-N2-C2
5	U	2	NAG	O5-C5-C6-O6
2	8	2	NAG	O5-C5-C6-O6
5	r	2	NAG	O5-C5-C6-O6
2	k	1	NAG	O5-C5-C6-O6
2	j	1	NAG	O7-C7-N2-C2
3	h	2	NAG	O7-C7-N2-C2
7	e	3	BMA	C4-C5-C6-O6
2	d	1	NAG	O7-C7-N2-C2
2	k	2	NAG	O7-C7-N2-C2
3	P	2	NAG	O7-C7-N2-C2
2	N	2	NAG	C8-C7-N2-C2
2	BA	1	NAG	C8-C7-N2-C2
2	c	2	NAG	O5-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
2	T	2	NAG	O5-C5-C6-O6
2	Z	1	NAG	C4-C5-C6-O6
8	W	1	NAG	C8-C7-N2-C2
2	u	2	NAG	O5-C5-C6-O6
4	i	2	NAG	O5-C5-C6-O6
3	M	1	NAG	C4-C5-C6-O6
2	g	2	NAG	O5-C5-C6-O6
2	0	2	NAG	O5-C5-C6-O6
2	x	3	BMA	O5-C5-C6-O6
2	AA	3	BMA	O5-C5-C6-O6
3	J	2	NAG	O5-C5-C6-O6
7	e	2	NAG	O5-C5-C6-O6
4	5	2	NAG	O5-C5-C6-O6
6	2	2	NAG	O5-C5-C6-O6
2	N	3	BMA	O5-C5-C6-O6
2	c	3	BMA	O5-C5-C6-O6
7	V	5	MAN	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	f	4	MAN	O5-C5-C6-O6
5	L	4	MAN	O5-C5-C6-O6
7	w	4	MAN	O5-C5-C6-O6
2	z	2	NAG	O5-C5-C6-O6
4	5	4	MAN	O5-C5-C6-O6
6	2	4	MAN	O5-C5-C6-O6
2	N	2	NAG	O7-C7-N2-C2
4	5	1	NAG	C8-C7-N2-C2
5	9	2	NAG	O5-C5-C6-O6
2	g	2	NAG	C3-C2-N2-C7
2	s	1	NAG	C3-C2-N2-C7
8	W	1	NAG	C3-C2-N2-C7
5	9	4	MAN	O5-C5-C6-O6
2	R	3	BMA	O5-C5-C6-O6
3	S	2	NAG	O5-C5-C6-O6
6	Q	5	MAN	O5-C5-C6-O6
3	EA	1	NAG	C8-C7-N2-C2
2	BA	1	NAG	O7-C7-N2-C2
8	W	1	NAG	O7-C7-N2-C2
5	U	2	NAG	C4-C5-C6-O6
3	P	1	NAG	C1-C2-N2-C7
8	W	1	NAG	C1-C2-N2-C7
2	g	2	NAG	C8-C7-N2-C2
3	M	1	NAG	O5-C5-C6-O6
4	5	1	NAG	O7-C7-N2-C2
6	2	3	BMA	C4-C5-C6-O6
2	g	2	NAG	O7-C7-N2-C2
3	EA	1	NAG	O7-C7-N2-C2
2	R	1	NAG	C3-C2-N2-C7
2	R	2	NAG	C3-C2-N2-C7
2	Y	2	NAG	C3-C2-N2-C7
2	4	2	NAG	C3-C2-N2-C7
3	O	1	NAG	C3-C2-N2-C7
3	O	2	NAG	C3-C2-N2-C7
3	P	1	NAG	C3-C2-N2-C7
4	5	2	NAG	C3-C2-N2-C7
5	U	2	NAG	C3-C2-N2-C7
5	X	1	NAG	C3-C2-N2-C7
5	X	2	NAG	C3-C2-N2-C7
5	m	1	NAG	C3-C2-N2-C7
5	m	2	NAG	C3-C2-N2-C7
5	3	1	NAG	C3-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
6	2	2	NAG	C3-C2-N2-C7
7	e	1	NAG	C3-C2-N2-C7
7	w	1	NAG	C3-C2-N2-C7
7	6	2	NAG	C3-C2-N2-C7
5	U	1	NAG	C4-C5-C6-O6
2	k	1	NAG	C4-C5-C6-O6
2	o	1	NAG	C8-C7-N2-C2
5	U	1	NAG	O5-C5-C6-O6
2	R	1	NAG	C1-C2-N2-C7
2	R	2	NAG	C1-C2-N2-C7
2	o	1	NAG	C1-C2-N2-C7
2	u	1	NAG	C1-C2-N2-C7
2	z	1	NAG	C1-C2-N2-C7
3	H	2	NAG	C1-C2-N2-C7
3	O	1	NAG	C1-C2-N2-C7
3	O	2	NAG	C1-C2-N2-C7
3	P	2	NAG	C1-C2-N2-C7
3	S	2	NAG	C1-C2-N2-C7
3	b	2	NAG	C1-C2-N2-C7
3	EA	2	NAG	C1-C2-N2-C7
4	p	1	NAG	C1-C2-N2-C7
4	5	1	NAG	C1-C2-N2-C7
4	5	2	NAG	C1-C2-N2-C7
5	U	2	NAG	C1-C2-N2-C7
5	3	1	NAG	C1-C2-N2-C7
7	e	1	NAG	C1-C2-N2-C7
10	l	6	MAN	O5-C5-C6-O6
3	v	1	NAG	O5-C5-C6-O6
3	v	1	NAG	C4-C5-C6-O6
2	1	2	NAG	O5-C5-C6-O6
2	o	1	NAG	O7-C7-N2-C2

There are no ring outliers.

33 monomers are involved in 28 short contacts:

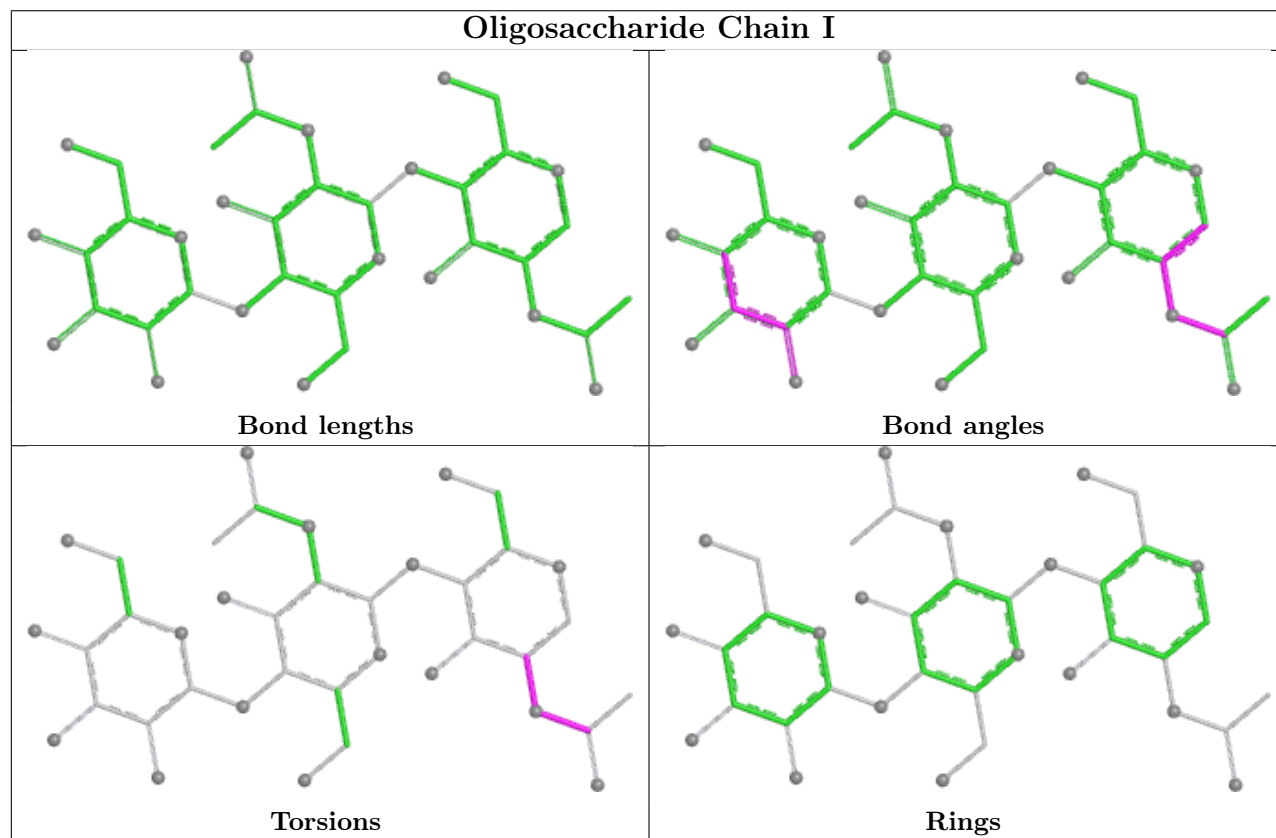
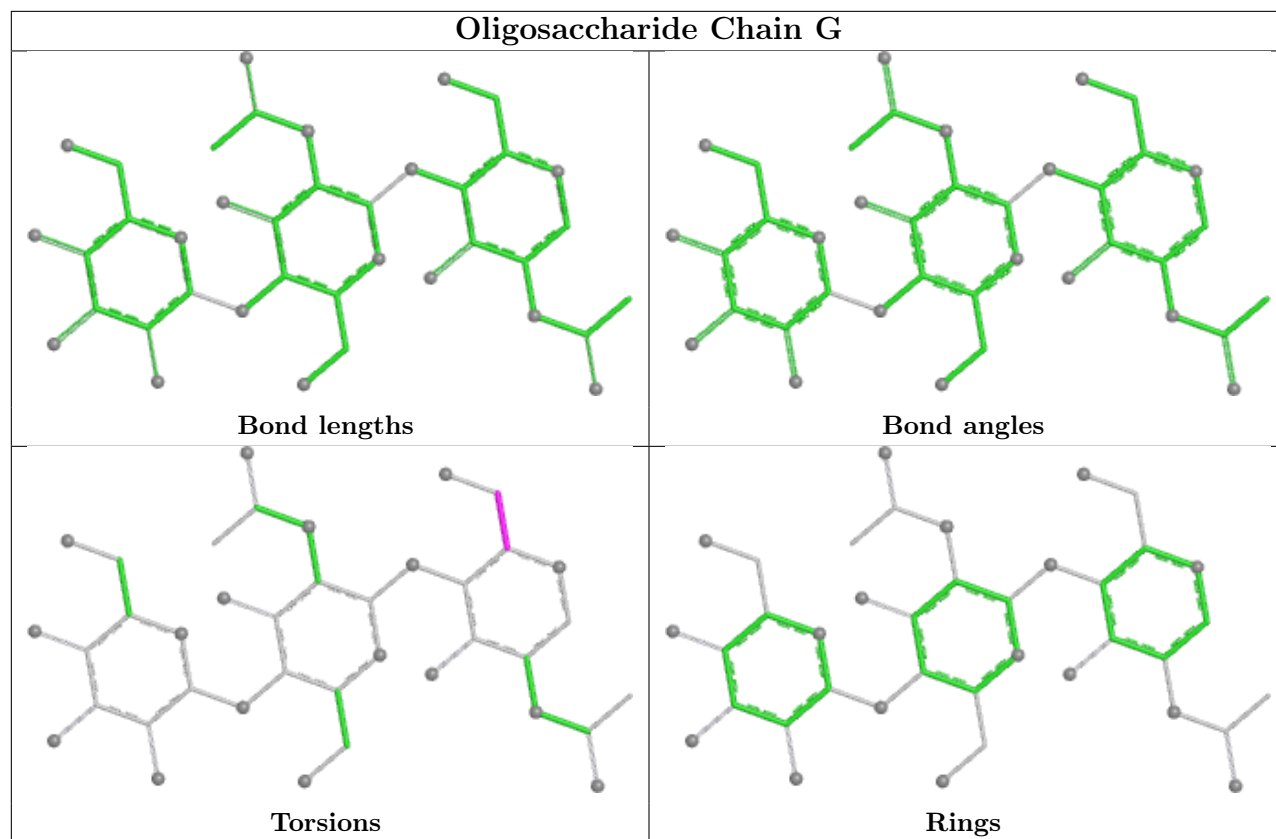
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	e	4	MAN	2	0
6	Q	1	NAG	1	0
11	7	1	NAG	1	0
3	O	1	NAG	1	0
10	l	3	BMA	1	0
7	V	3	BMA	3	0

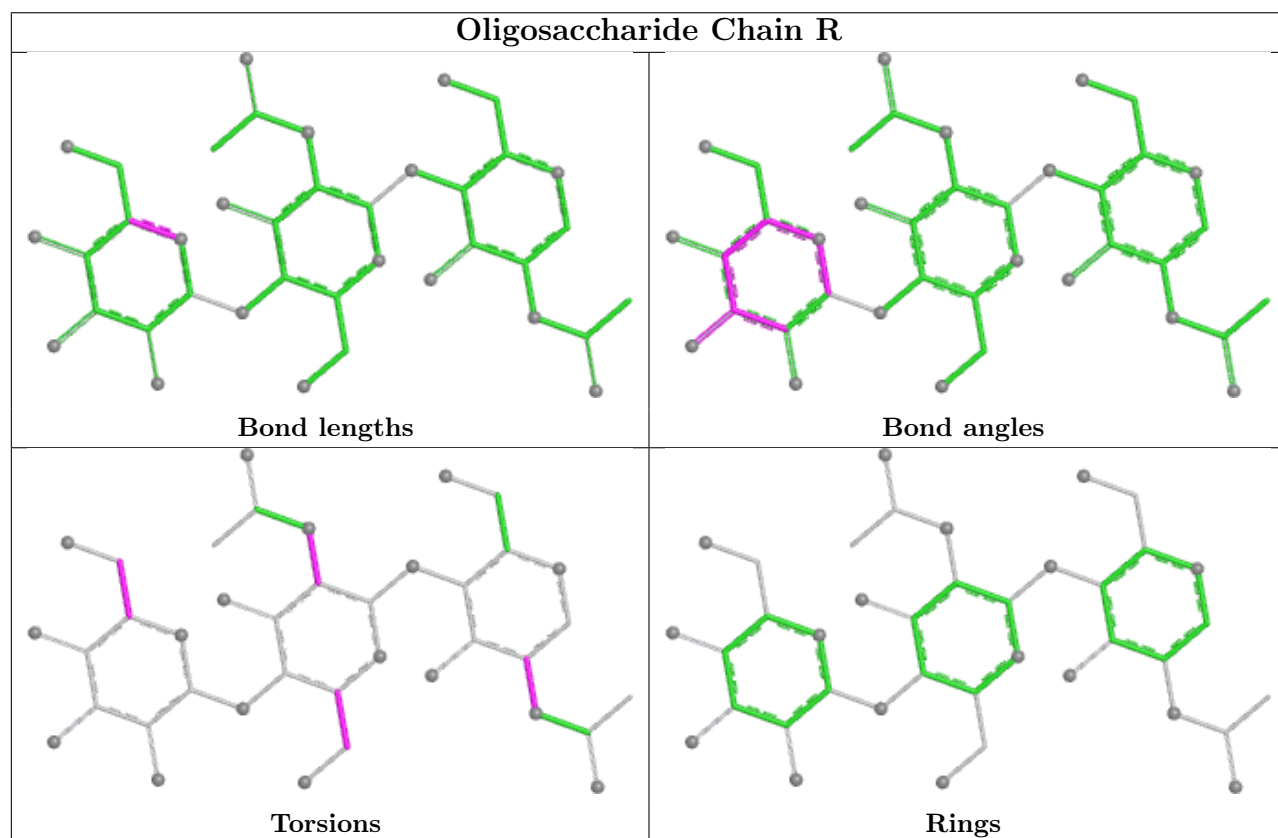
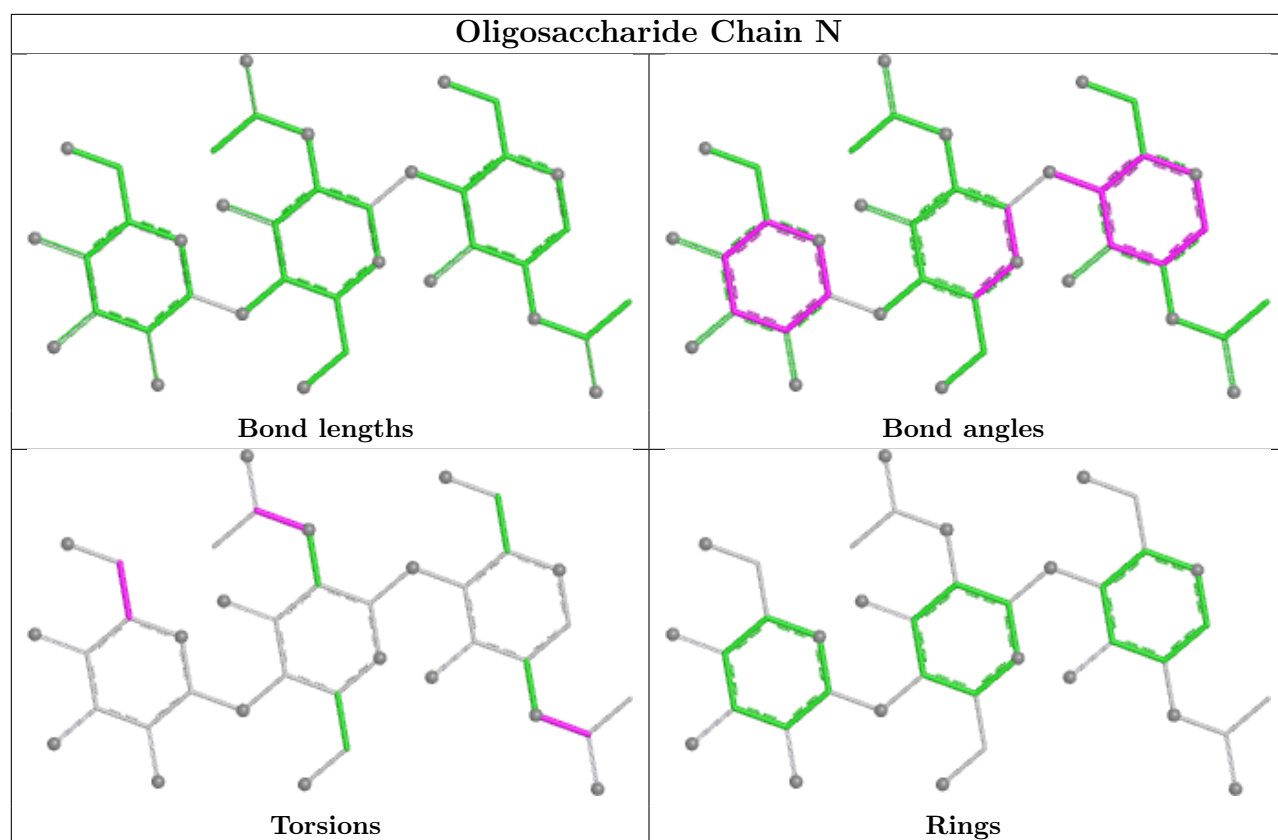
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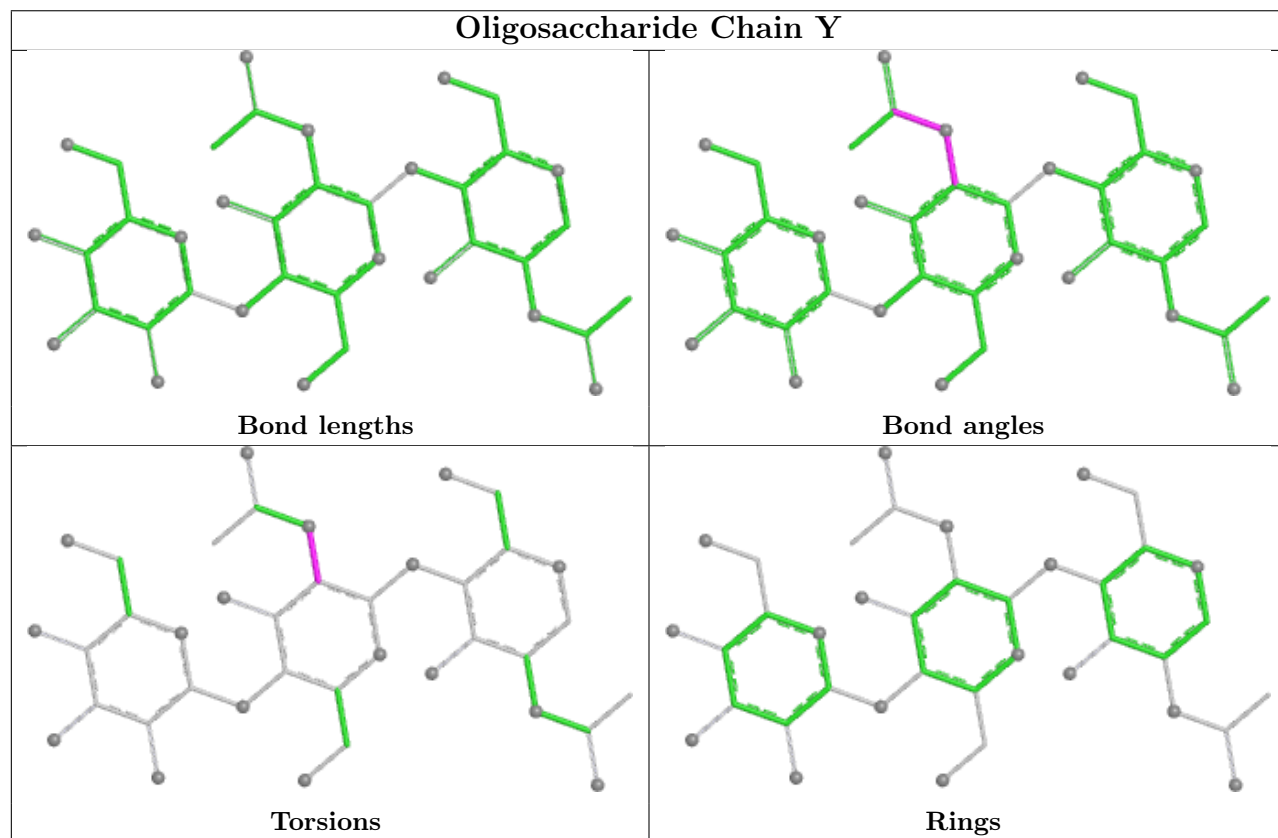
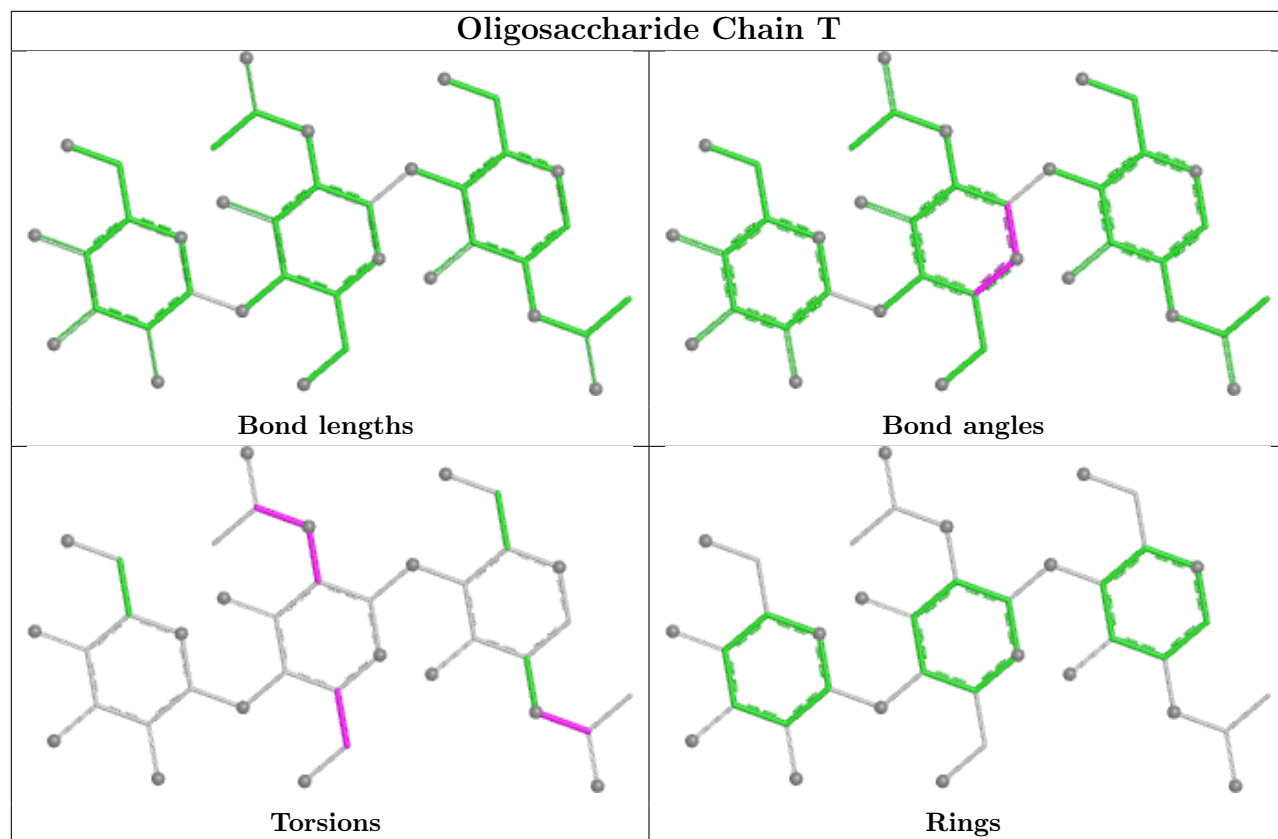
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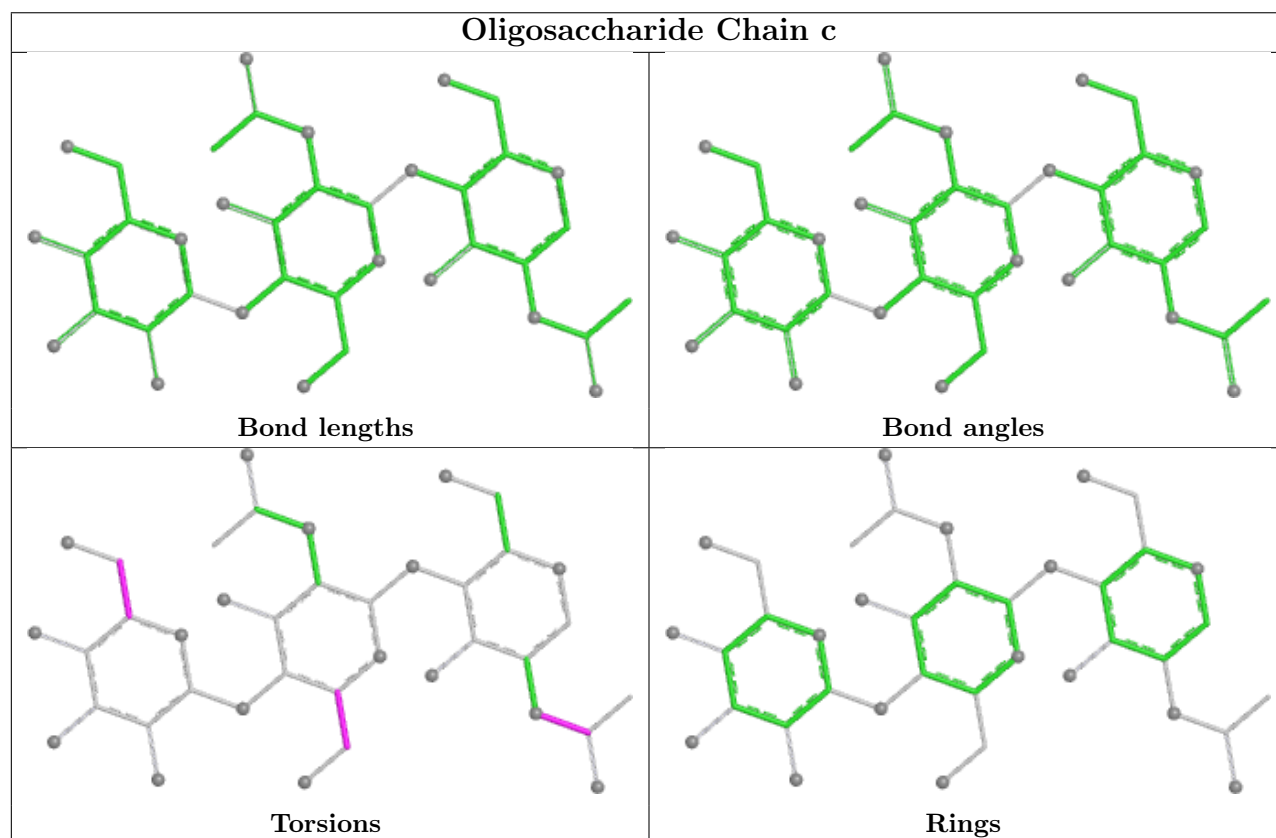
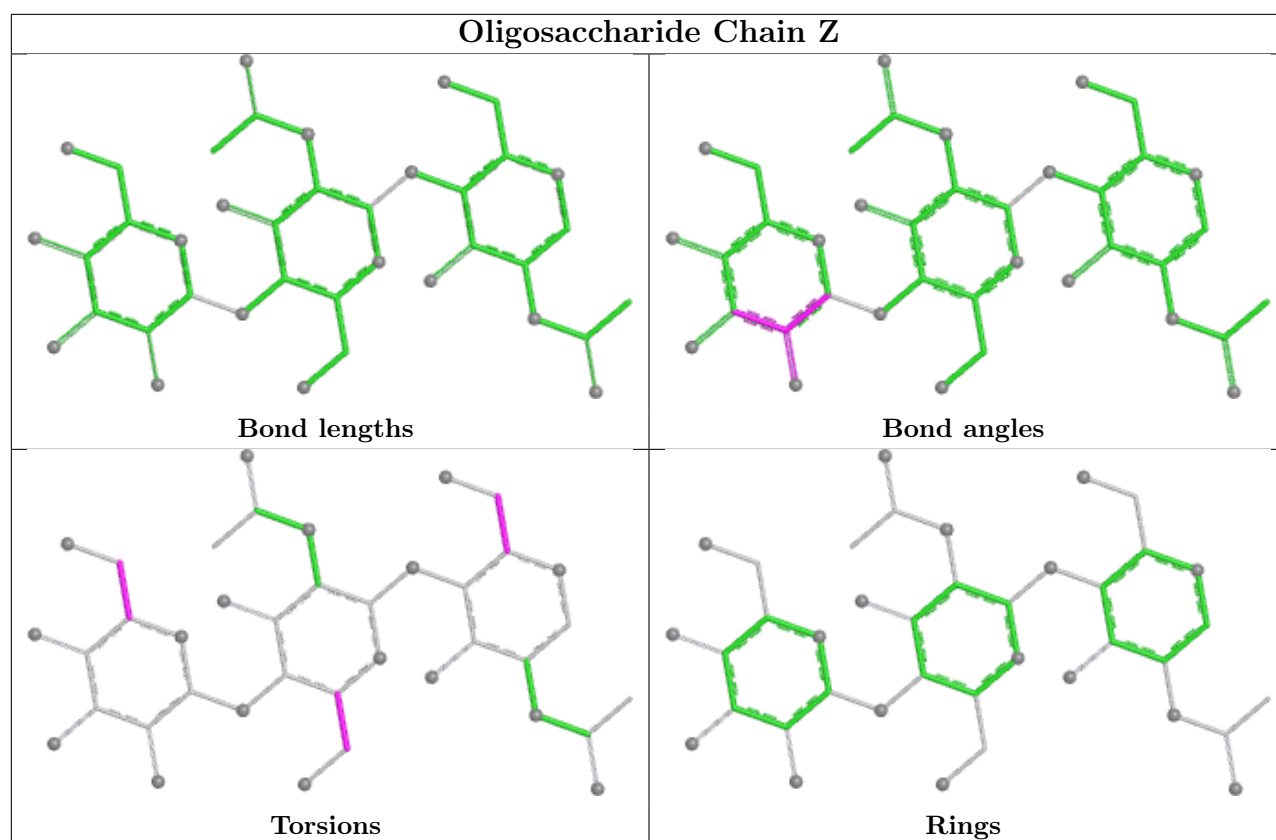
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	m	3	BMA	2	0
10	l	6	MAN	1	0
6	2	3	BMA	1	0
11	7	2	NAG	1	0
2	1	2	NAG	1	0
11	q	5	MAN	1	0
7	V	5	MAN	2	0
11	q	4	MAN	3	0
2	y	1	NAG	1	0
11	7	5	MAN	1	0
11	7	3	BMA	2	0
2	1	1	NAG	1	0
7	V	4	MAN	1	0
11	q	2	NAG	1	0
4	5	4	MAN	1	0
11	q	6	MAN	1	0
11	7	4	MAN	3	0
11	q	1	NAG	2	0
5	m	4	MAN	2	0
2	y	2	NAG	1	0
4	5	3	BMA	1	0
7	e	3	BMA	2	0
3	P	1	NAG	1	0
11	q	3	BMA	2	0
6	2	2	NAG	1	0
5	9	1	NAG	2	0
6	2	6	MAN	1	0

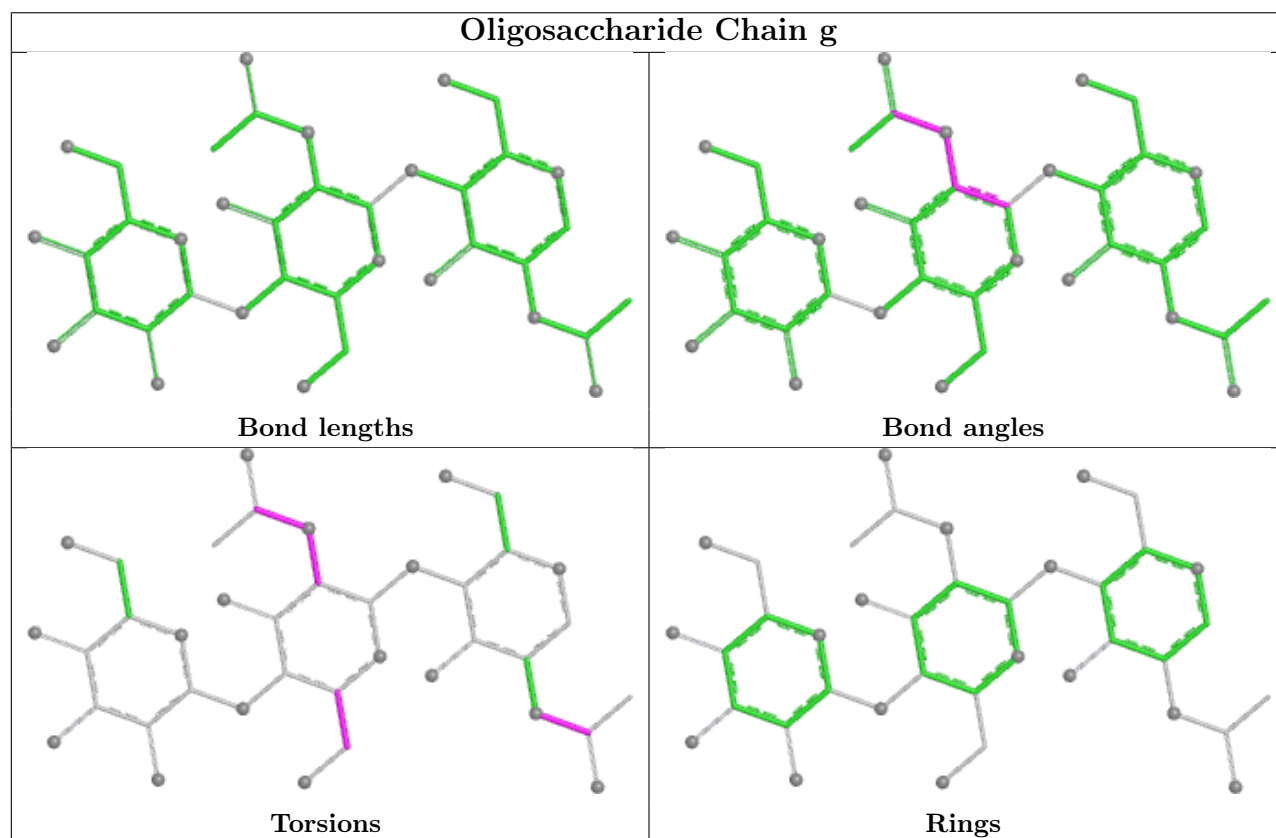
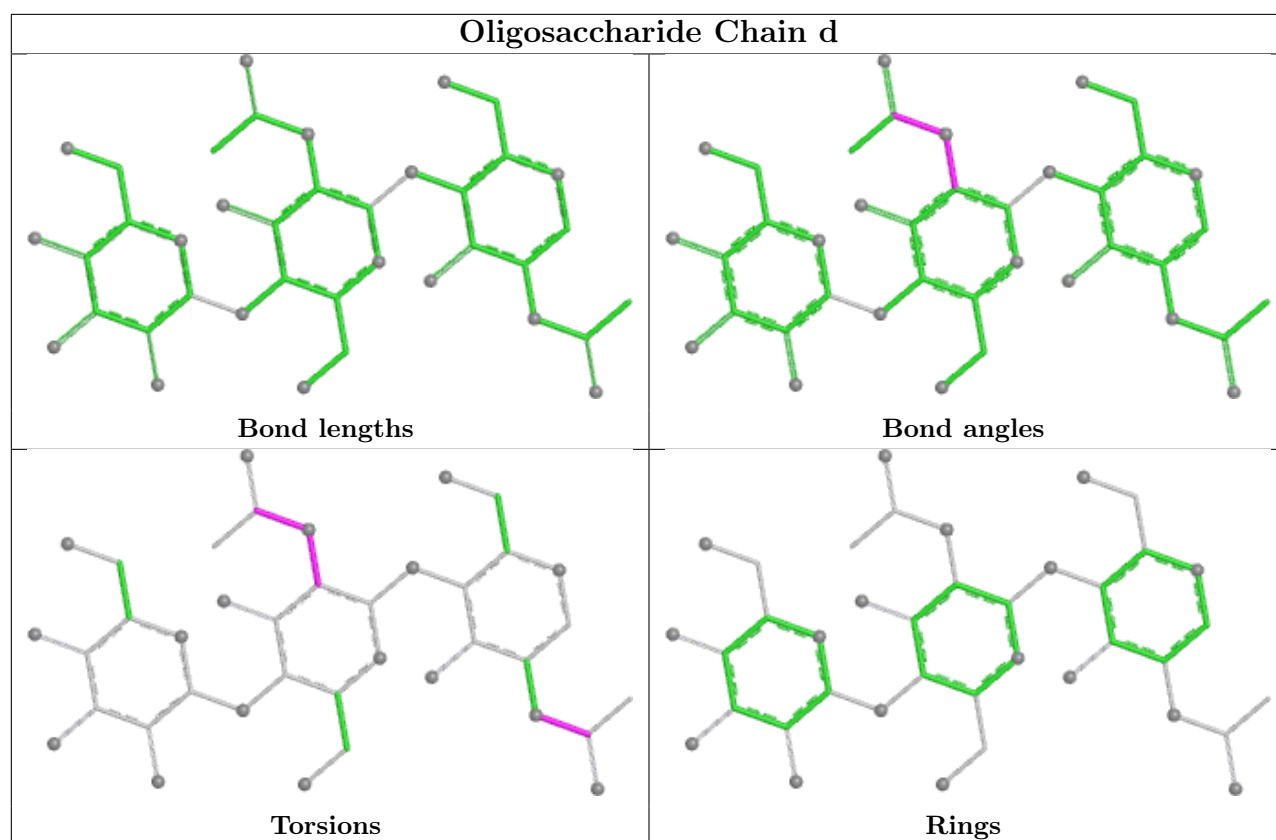
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

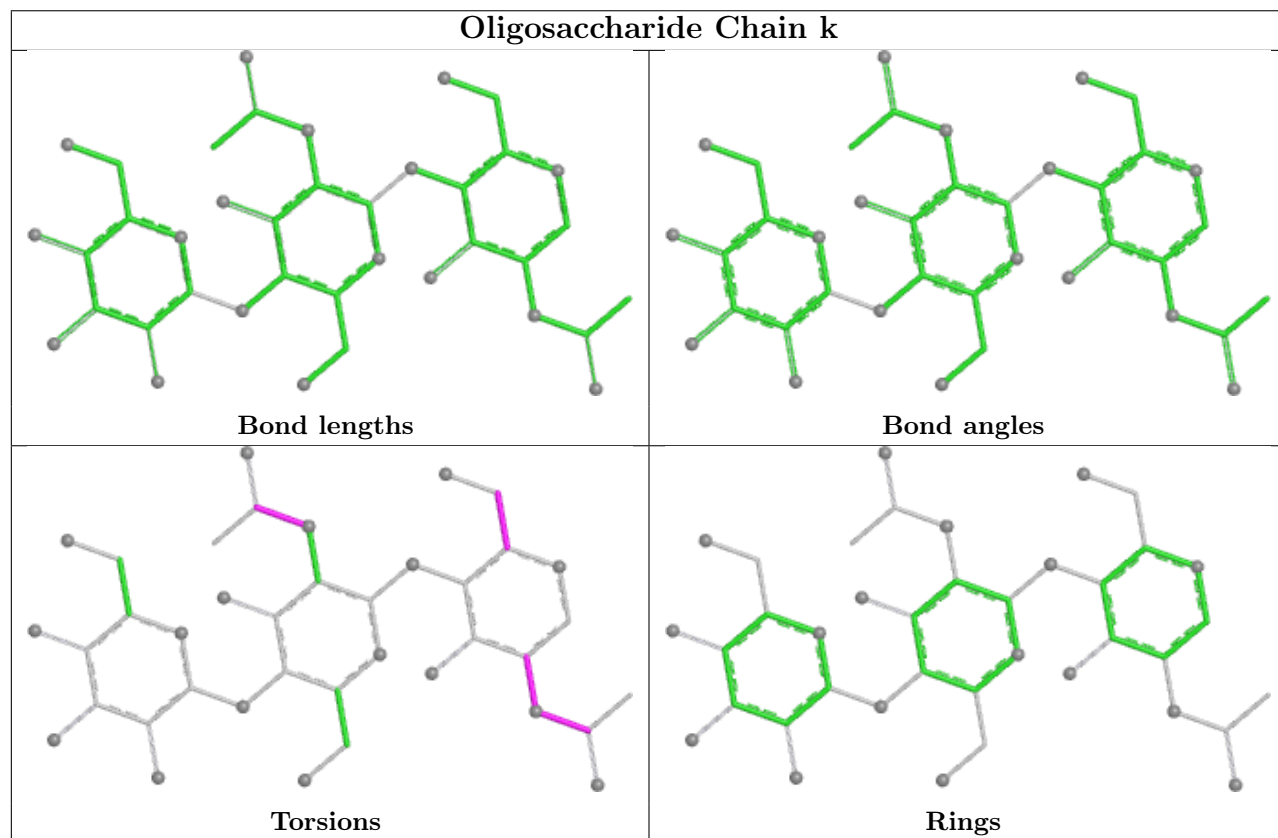
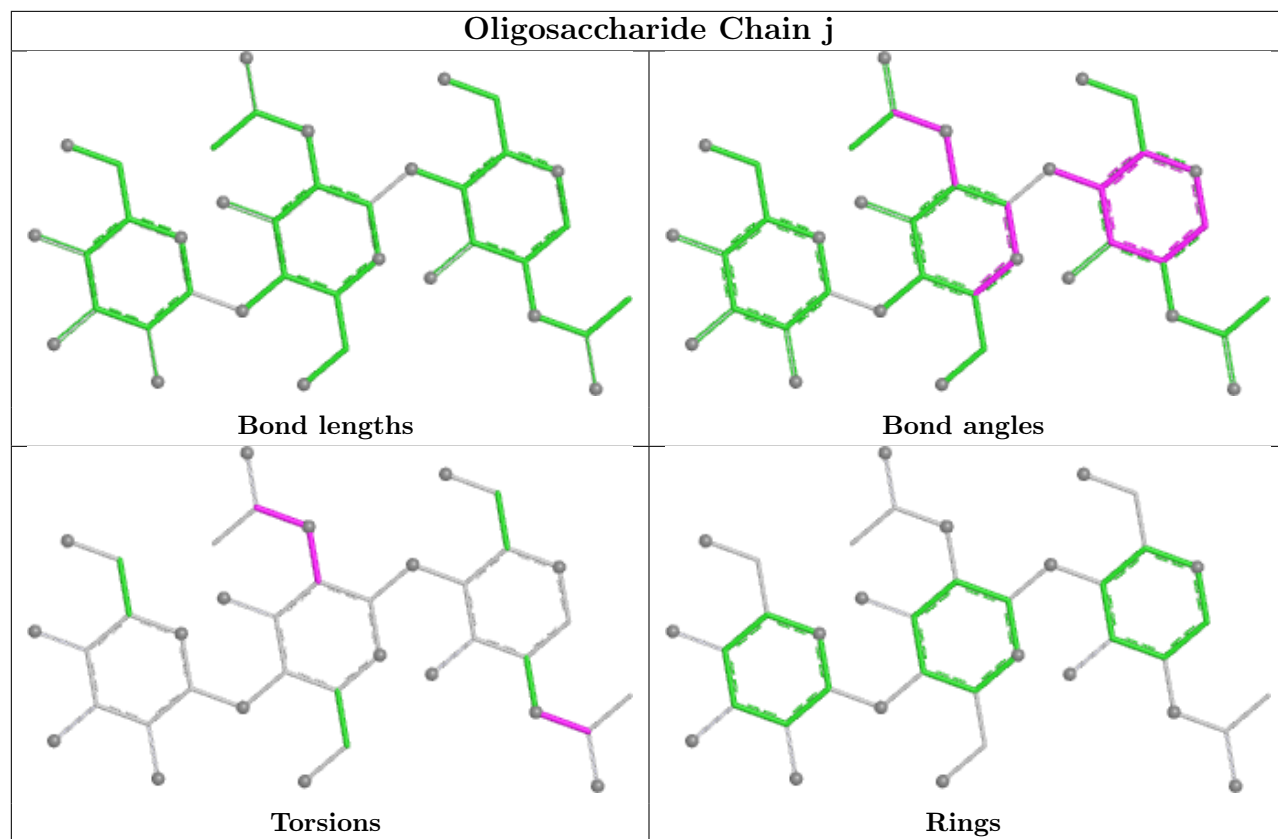


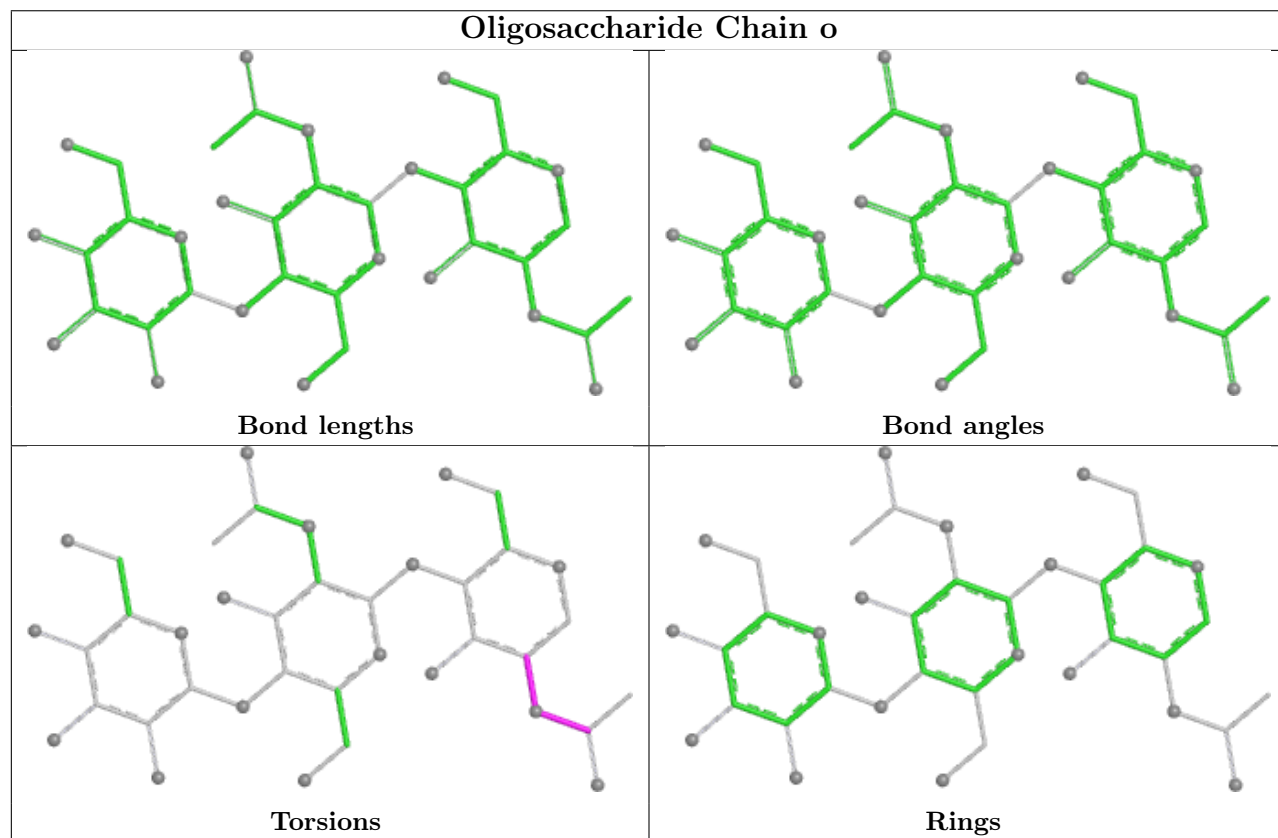
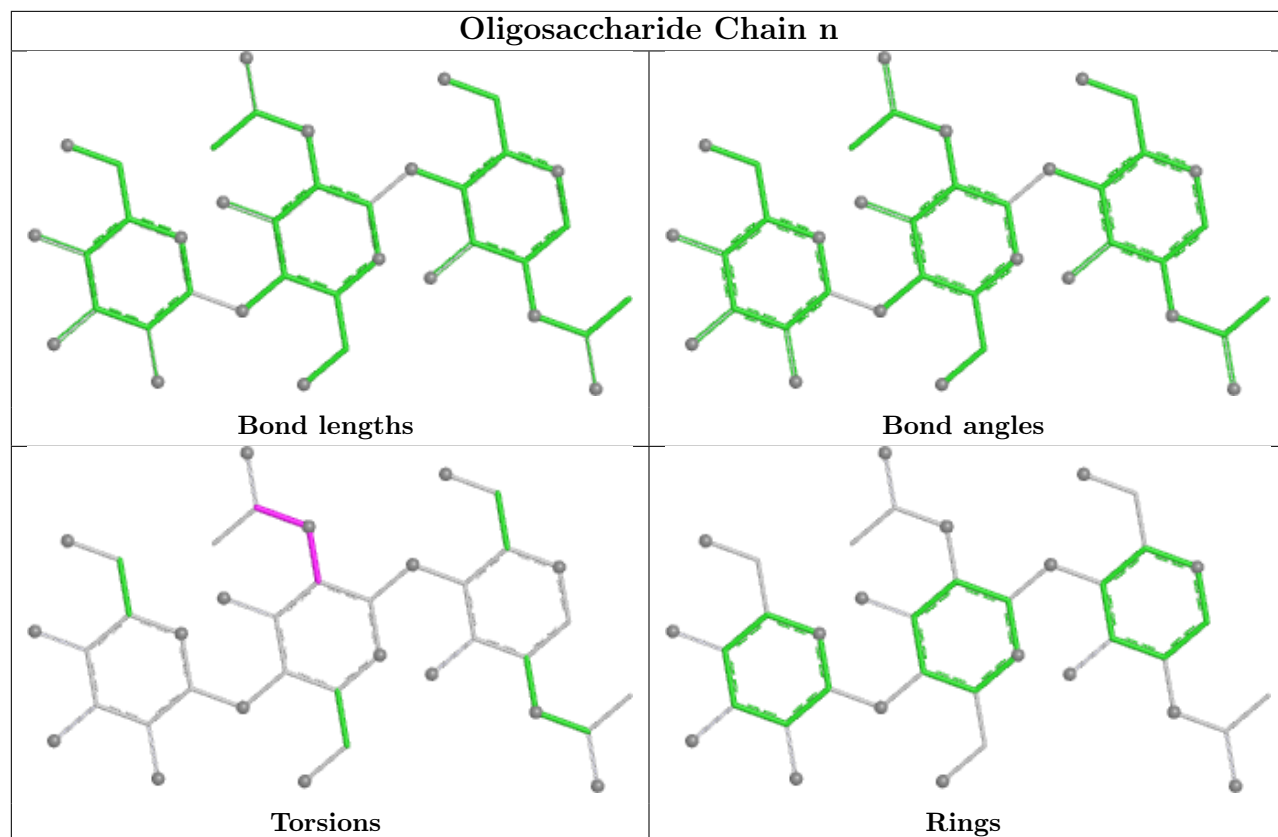


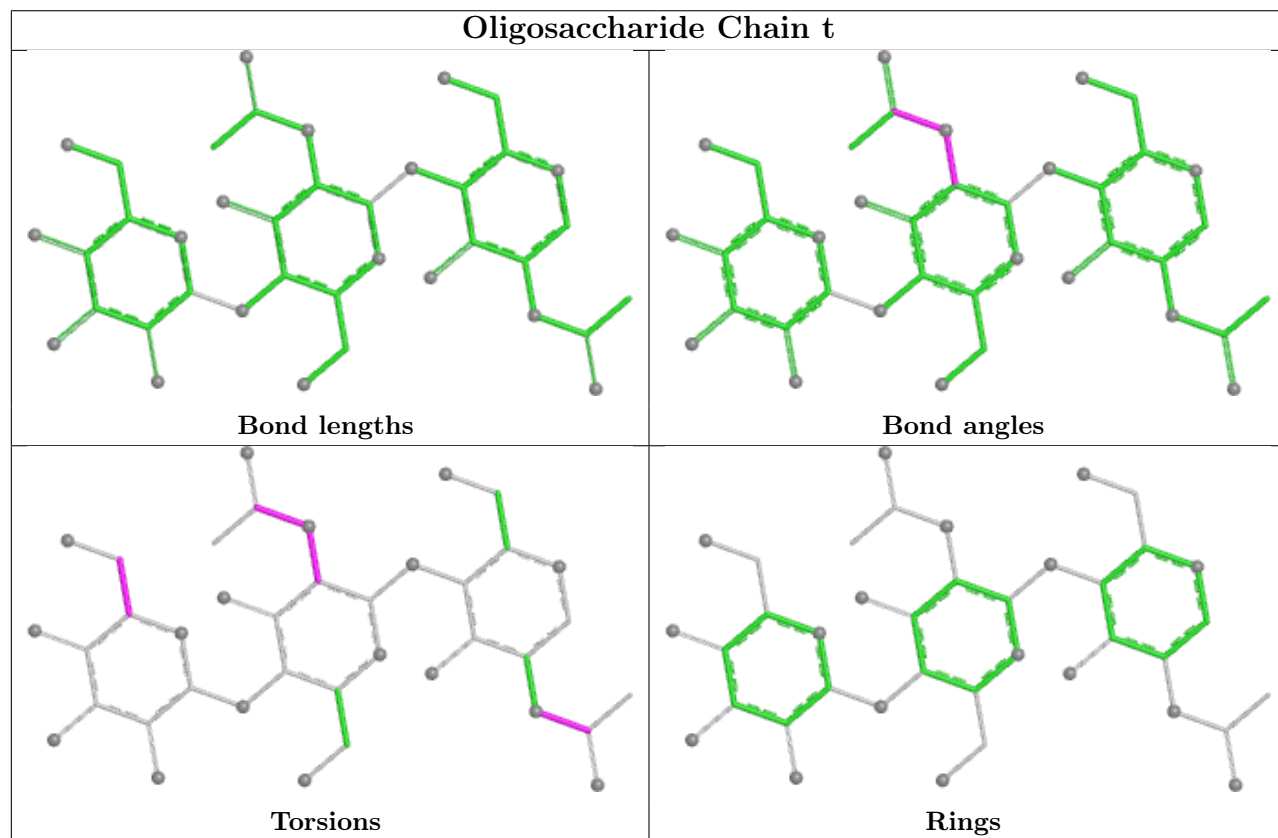
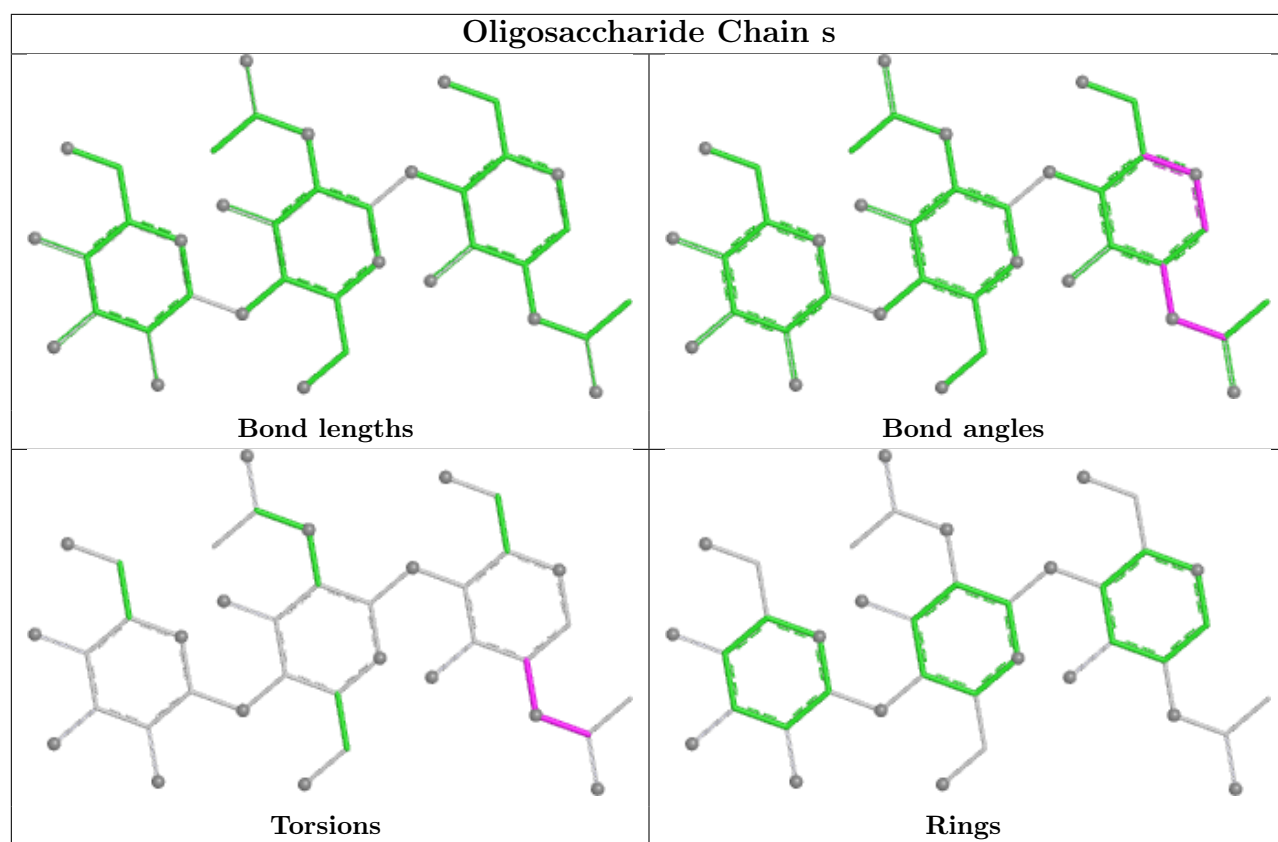


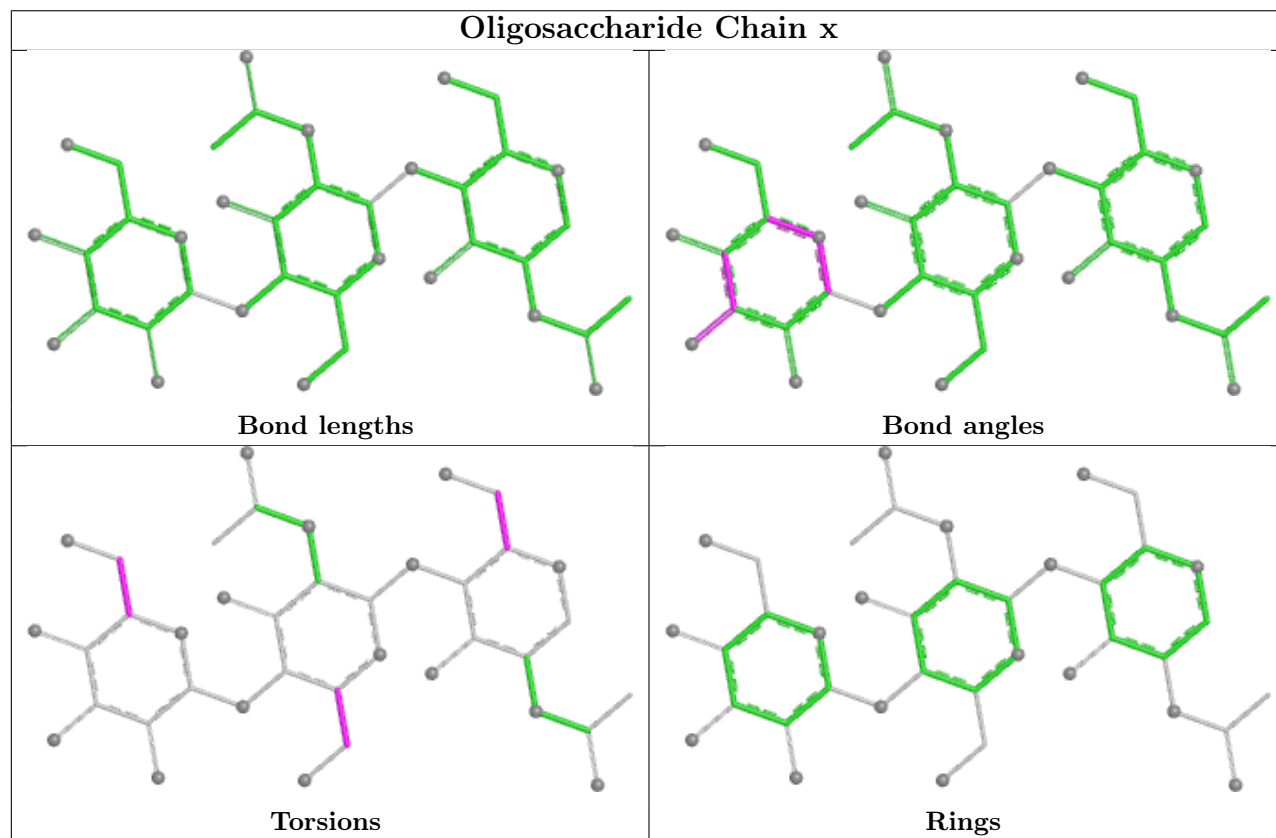
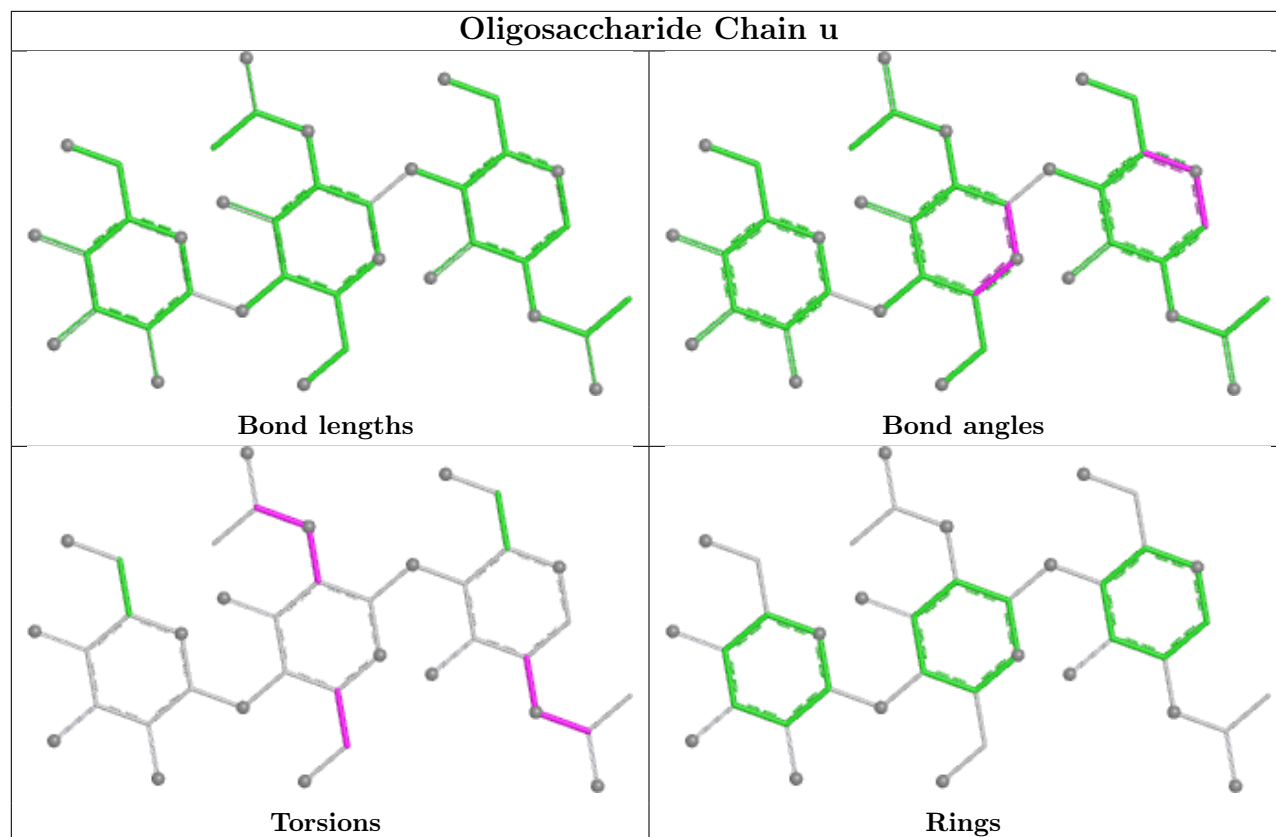


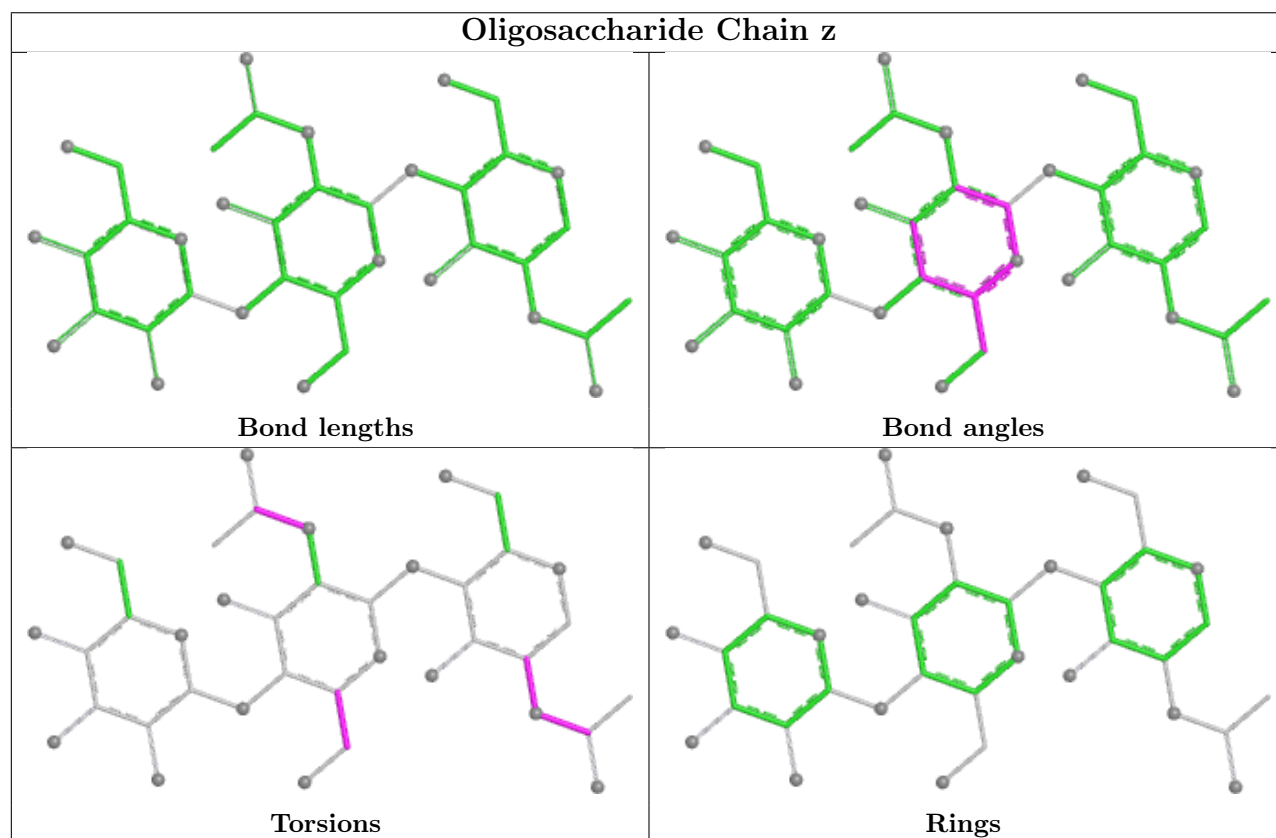
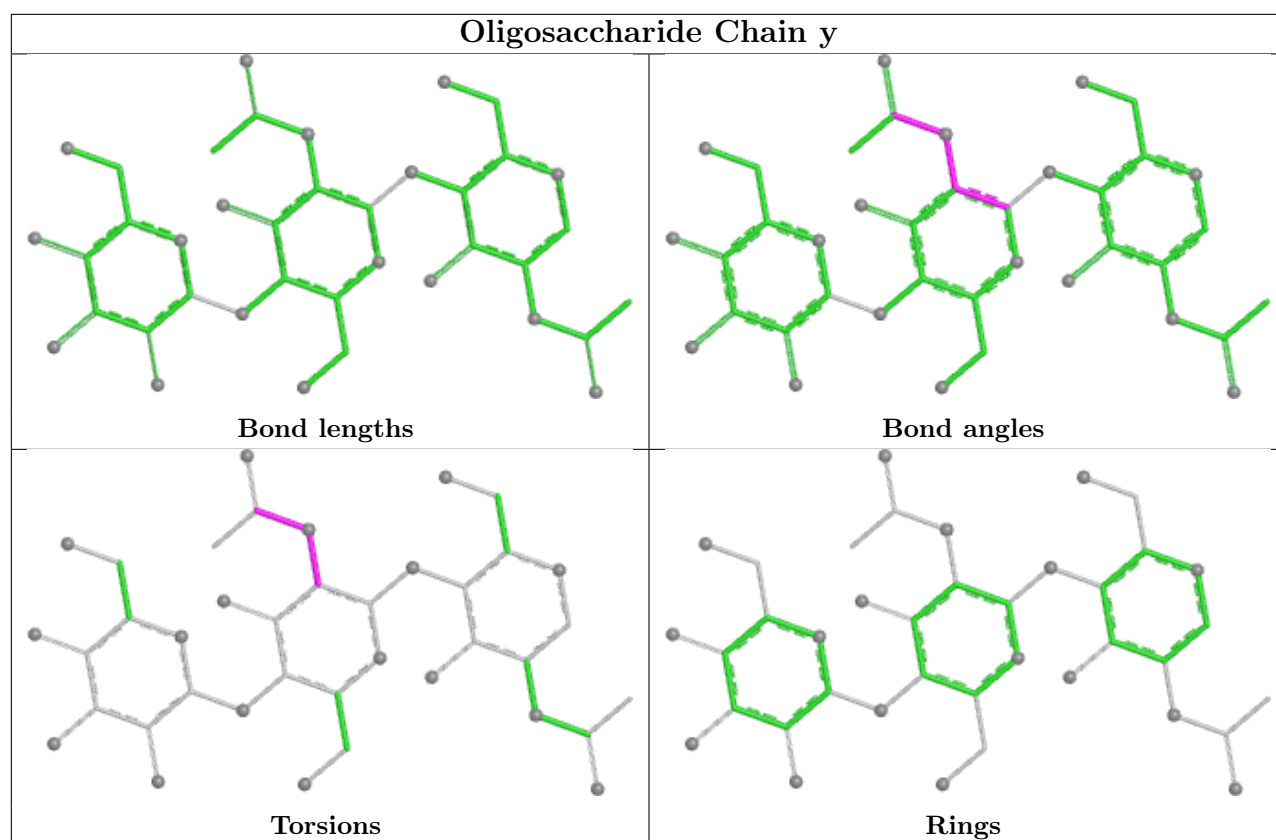


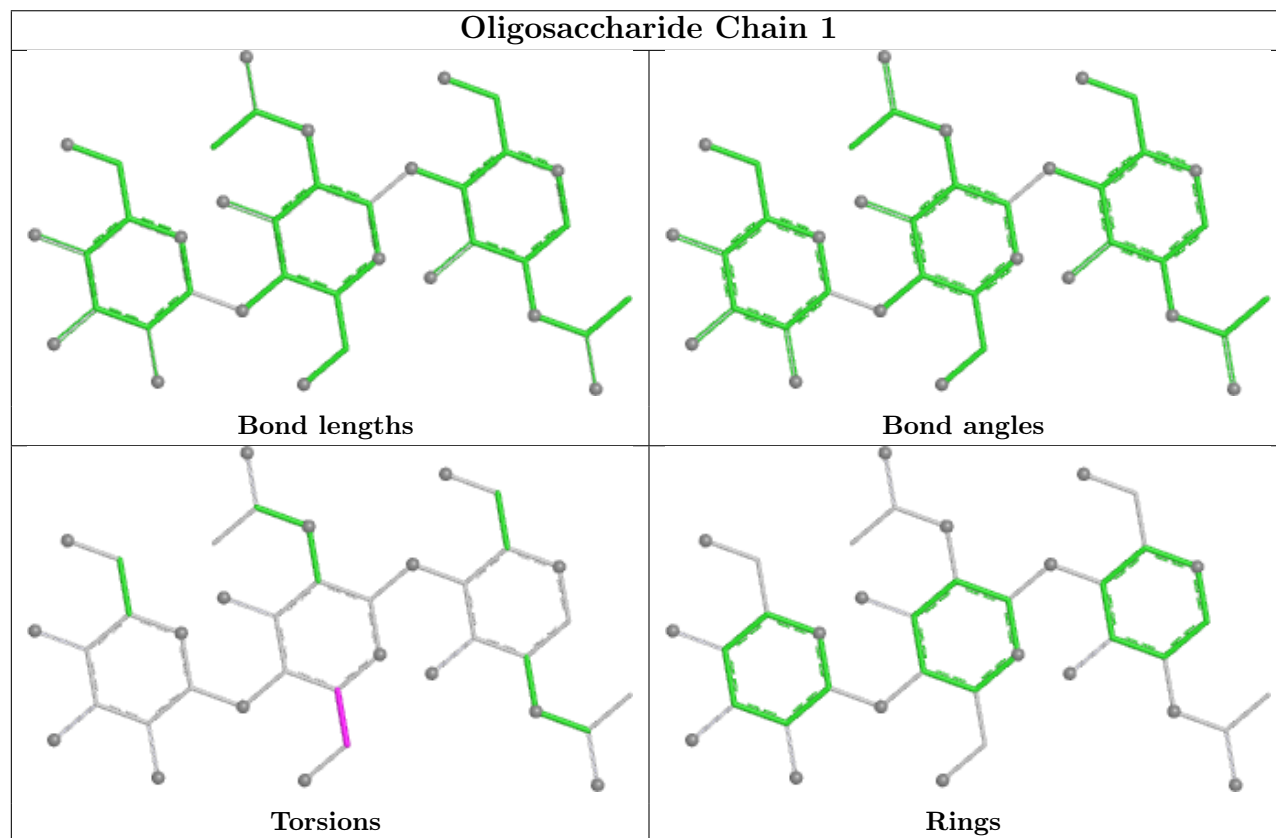
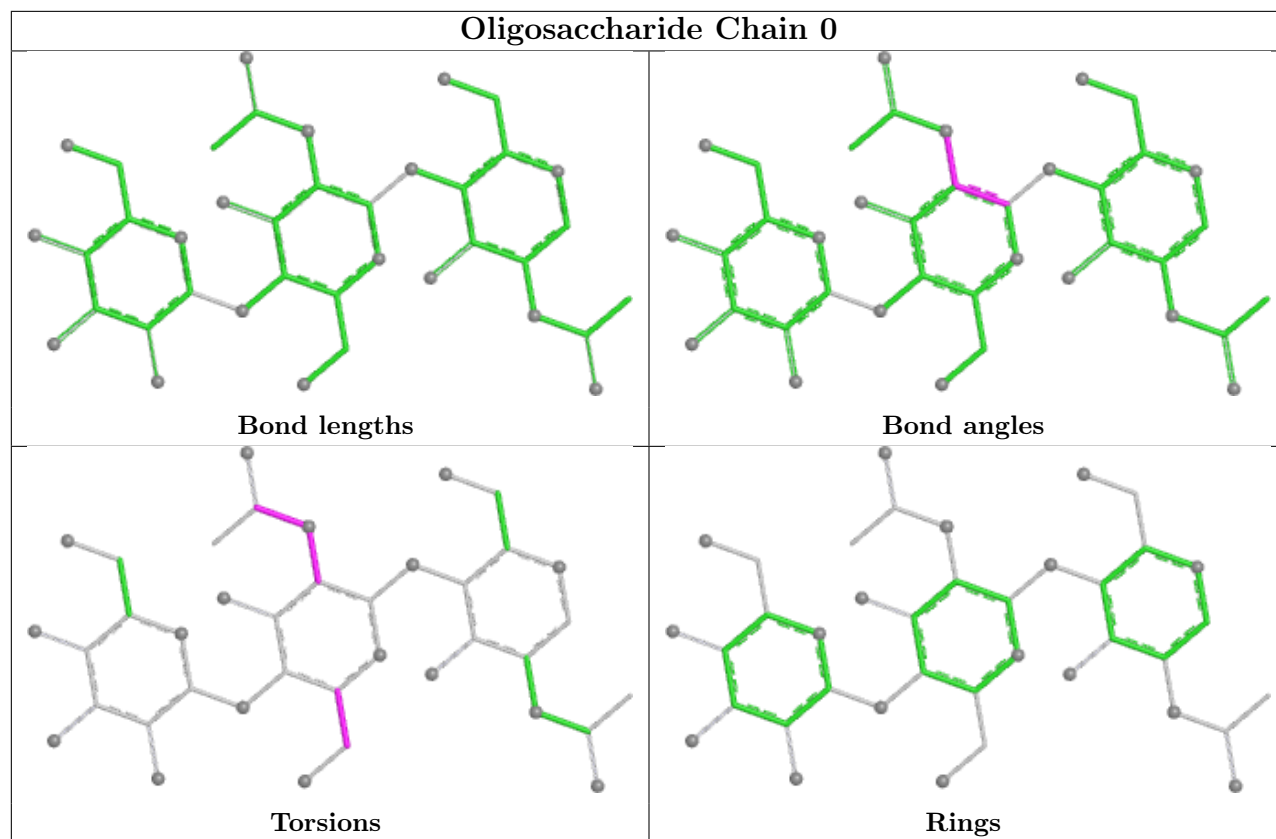


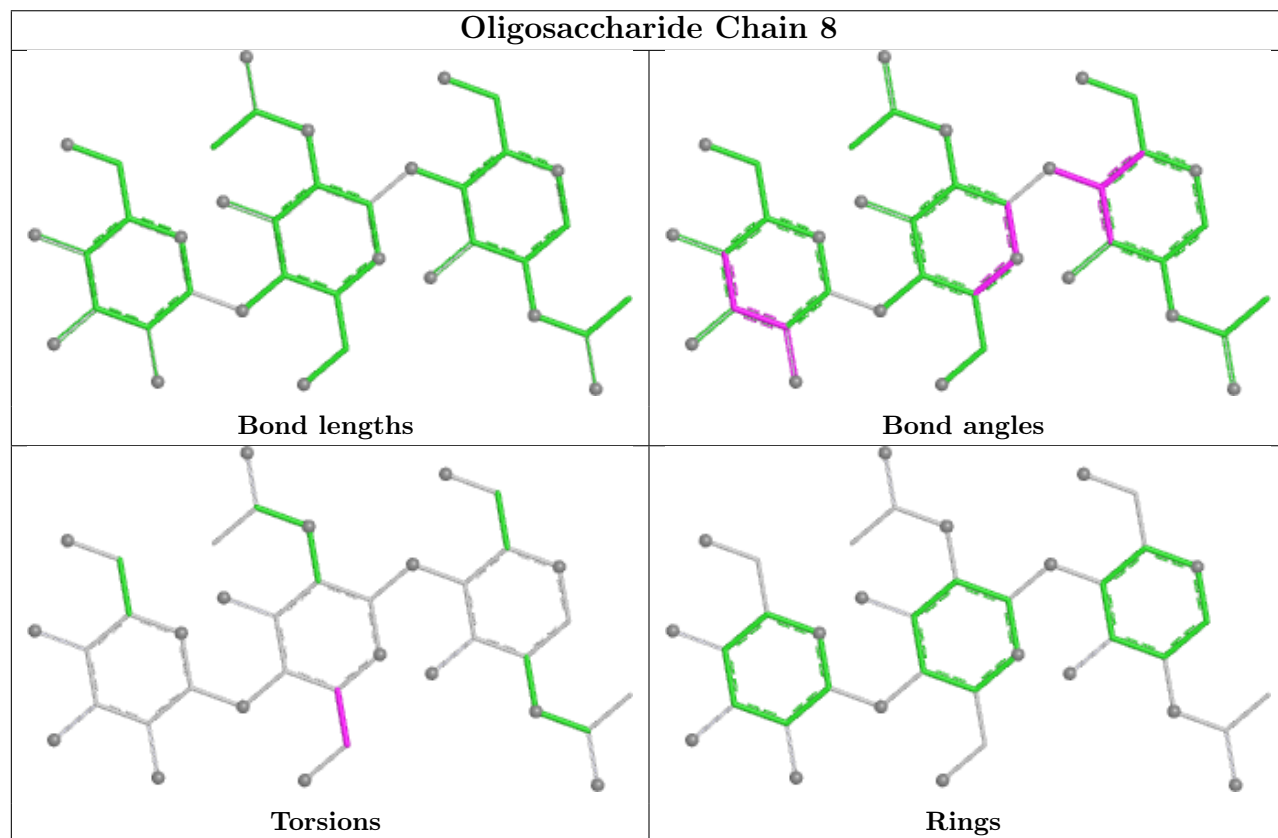
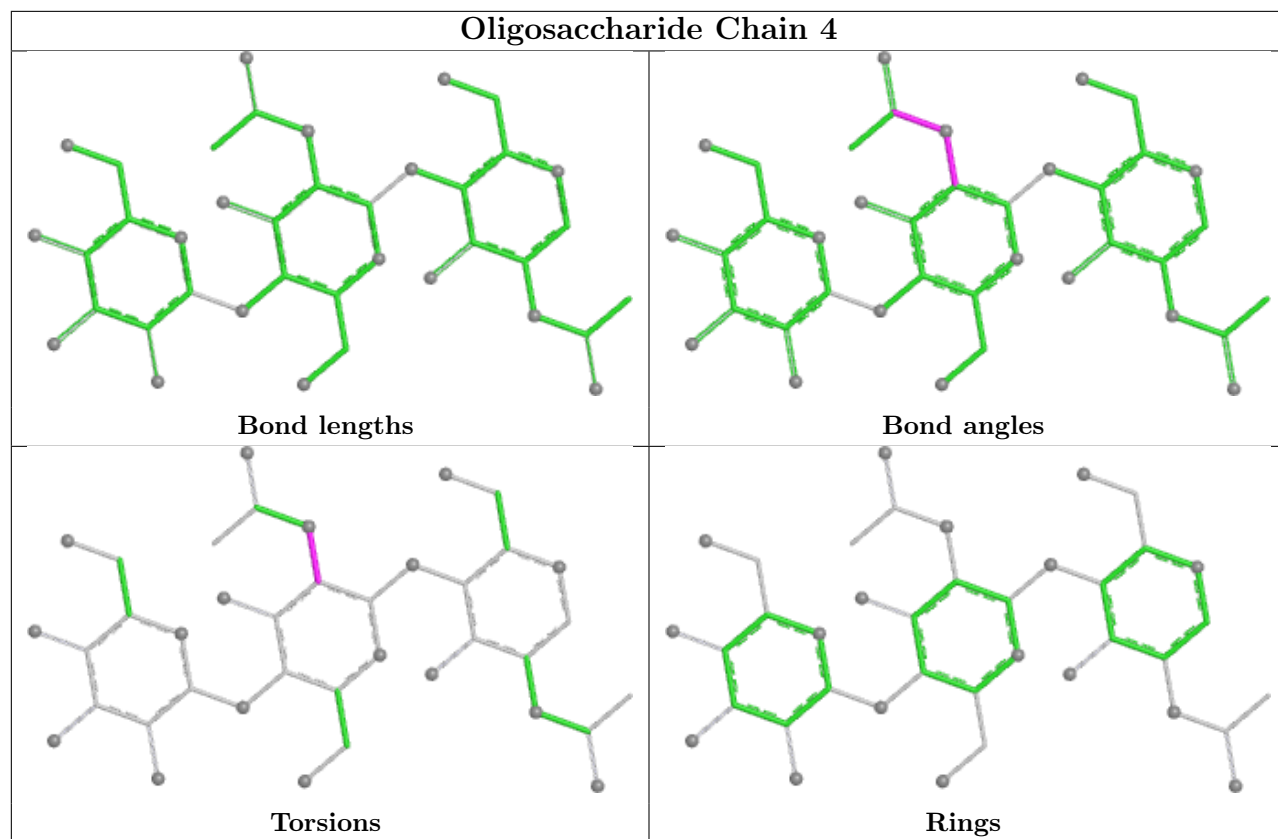


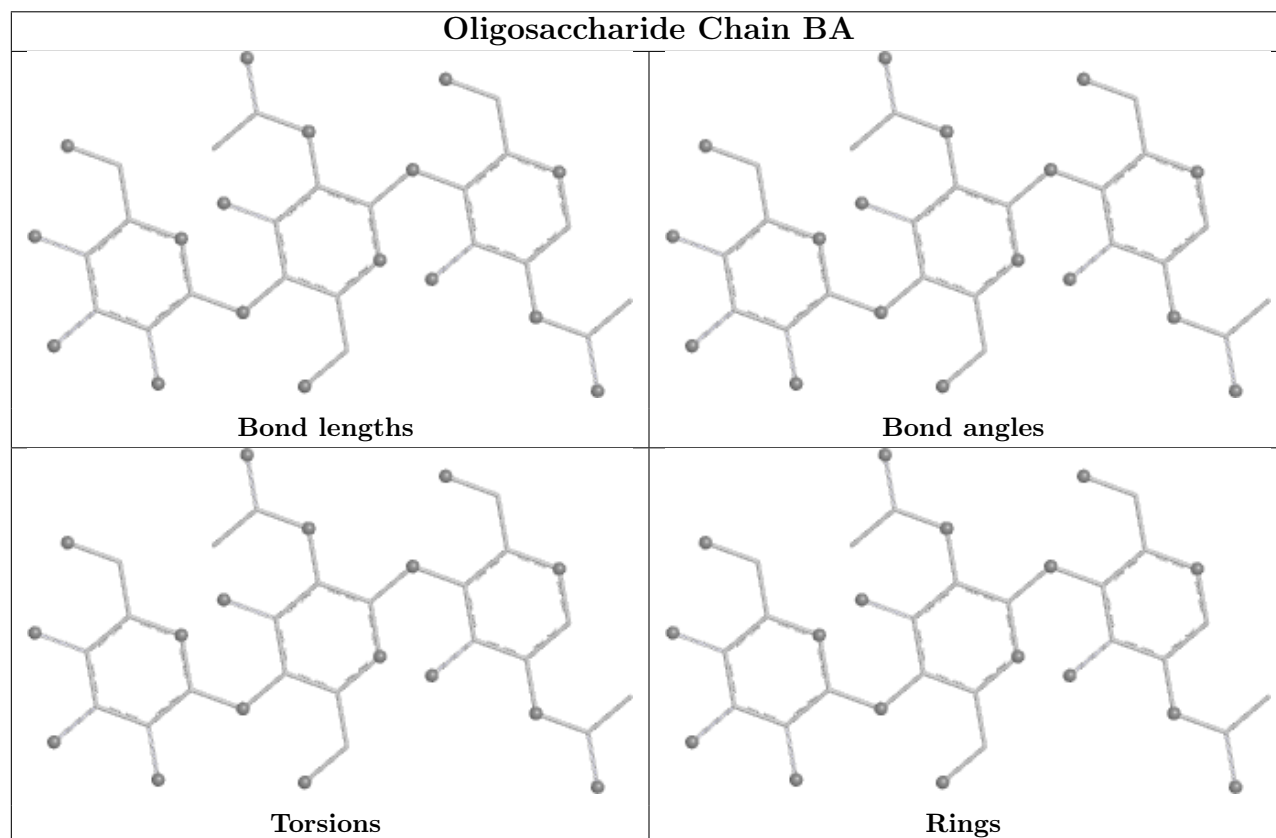
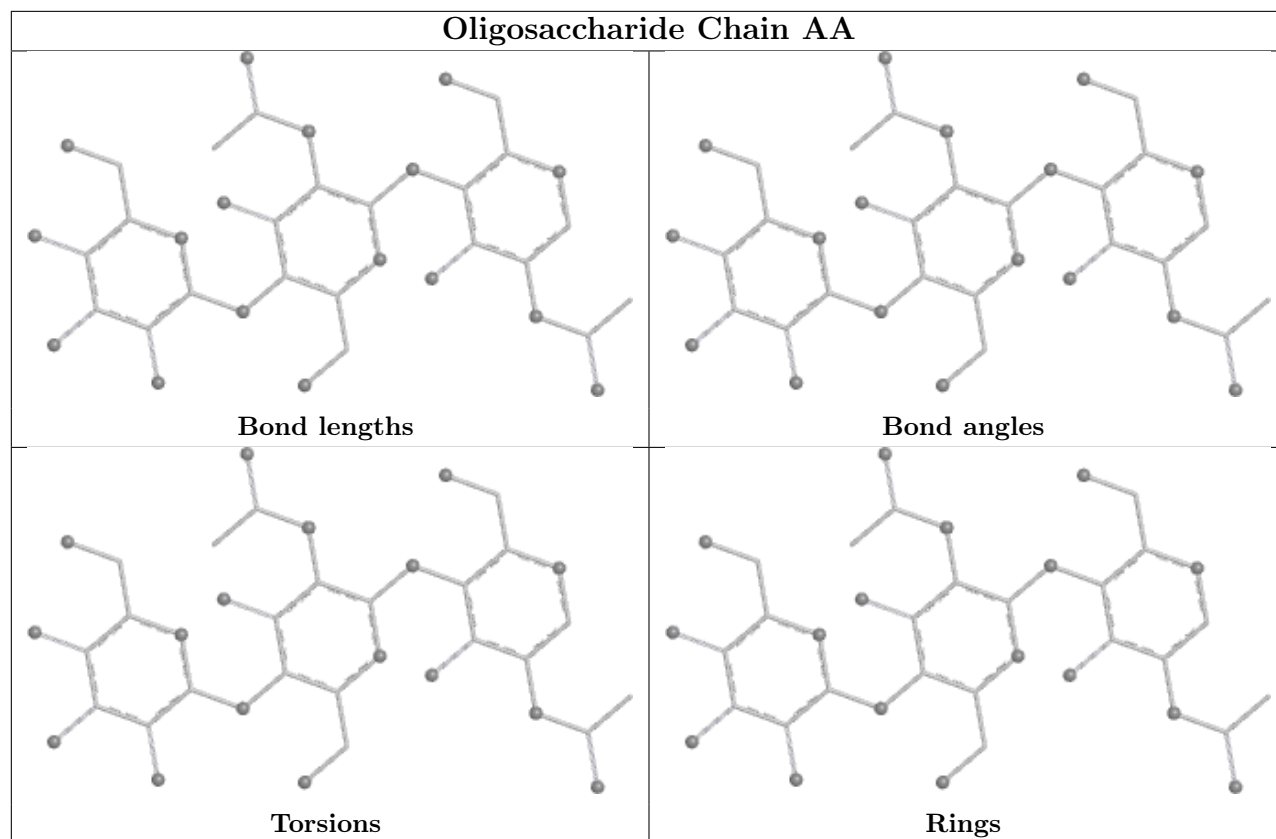


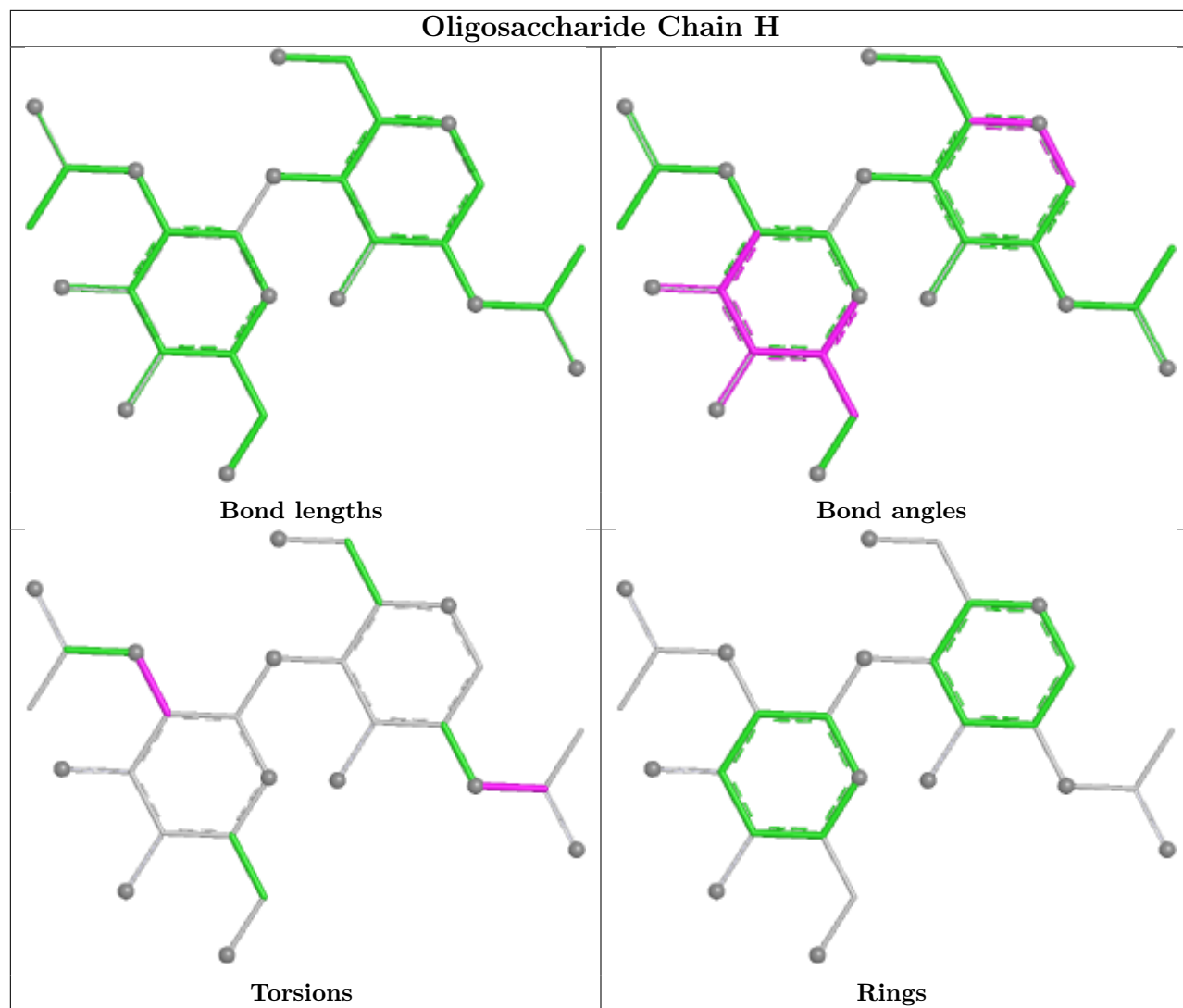


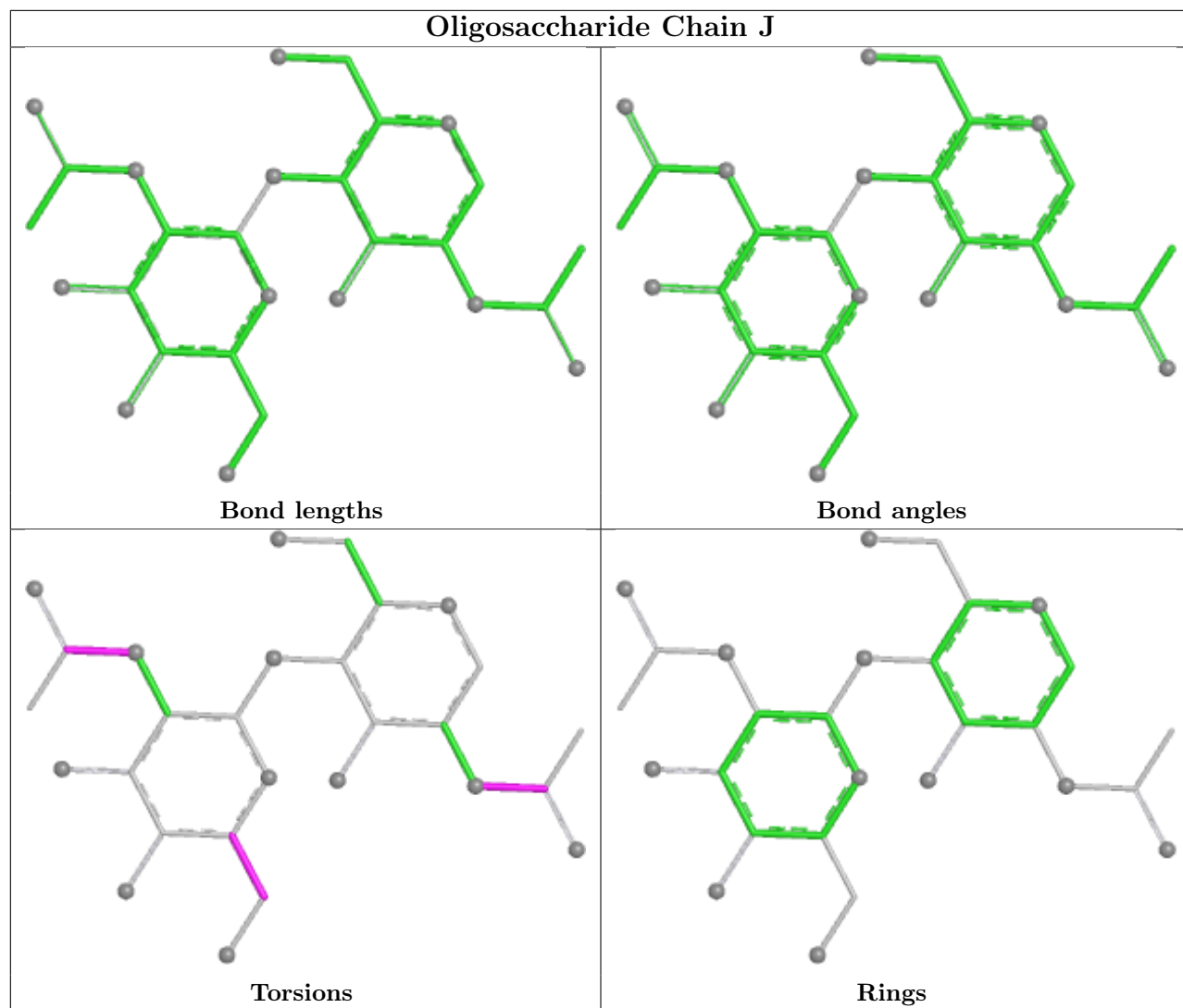


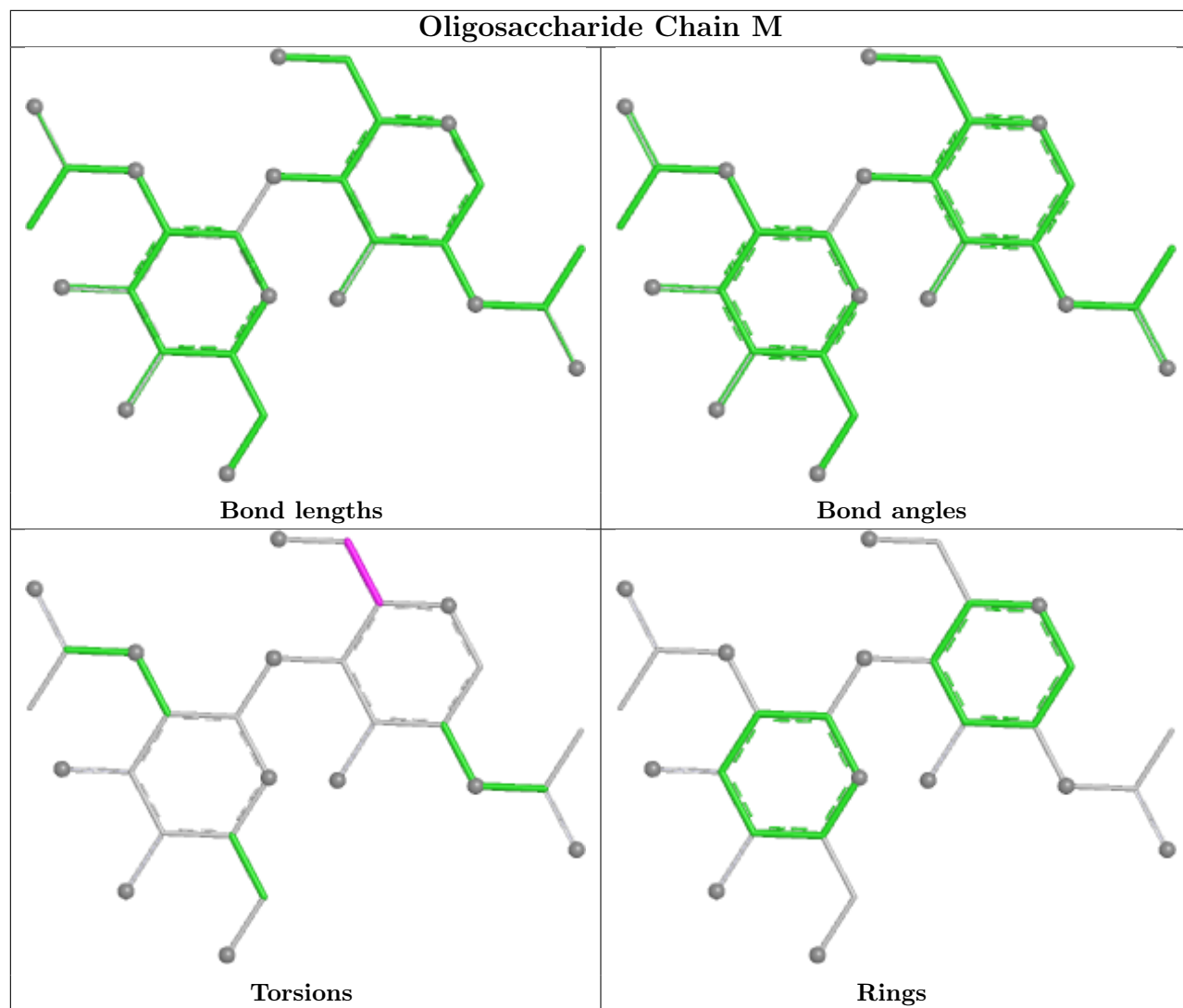


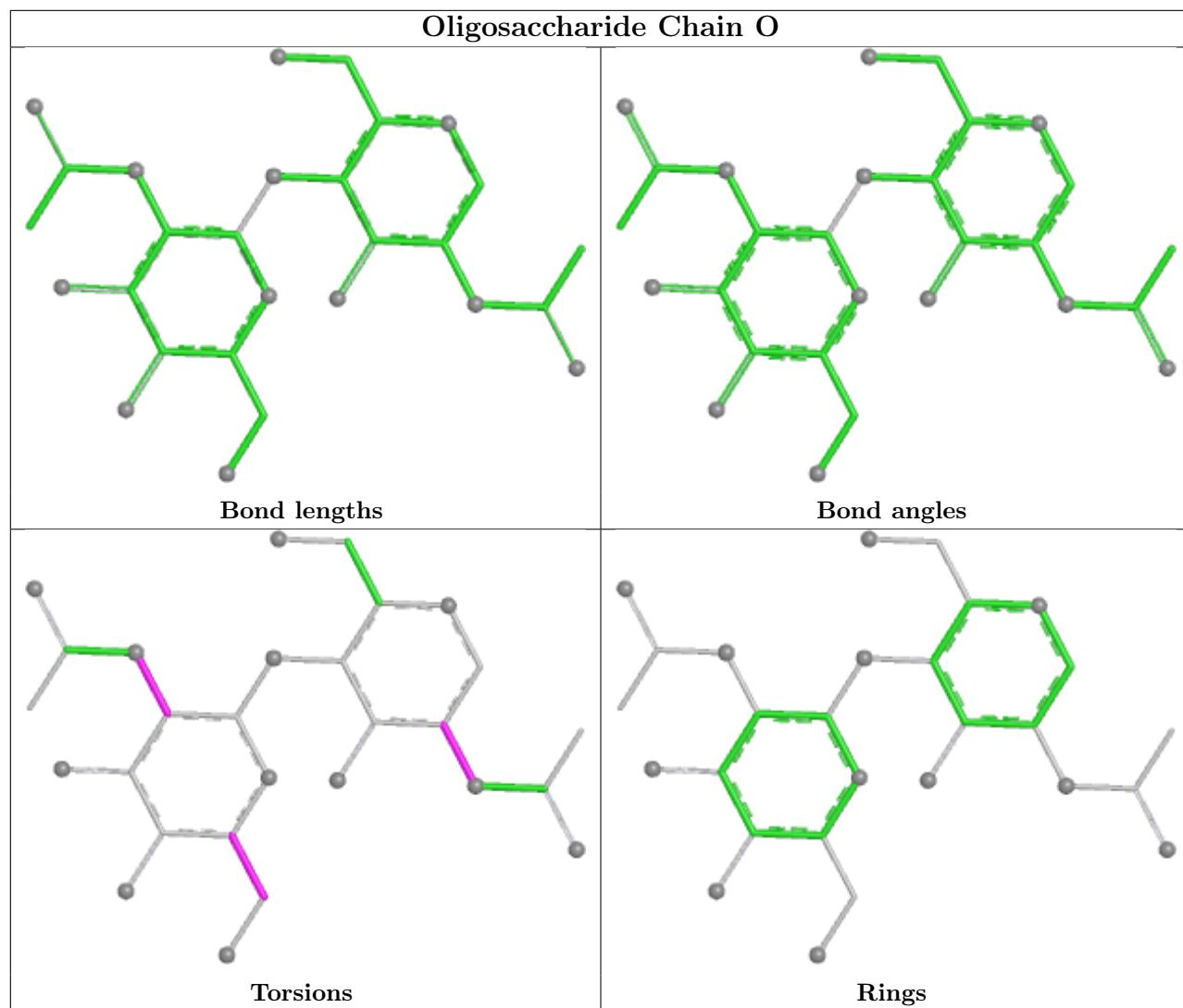


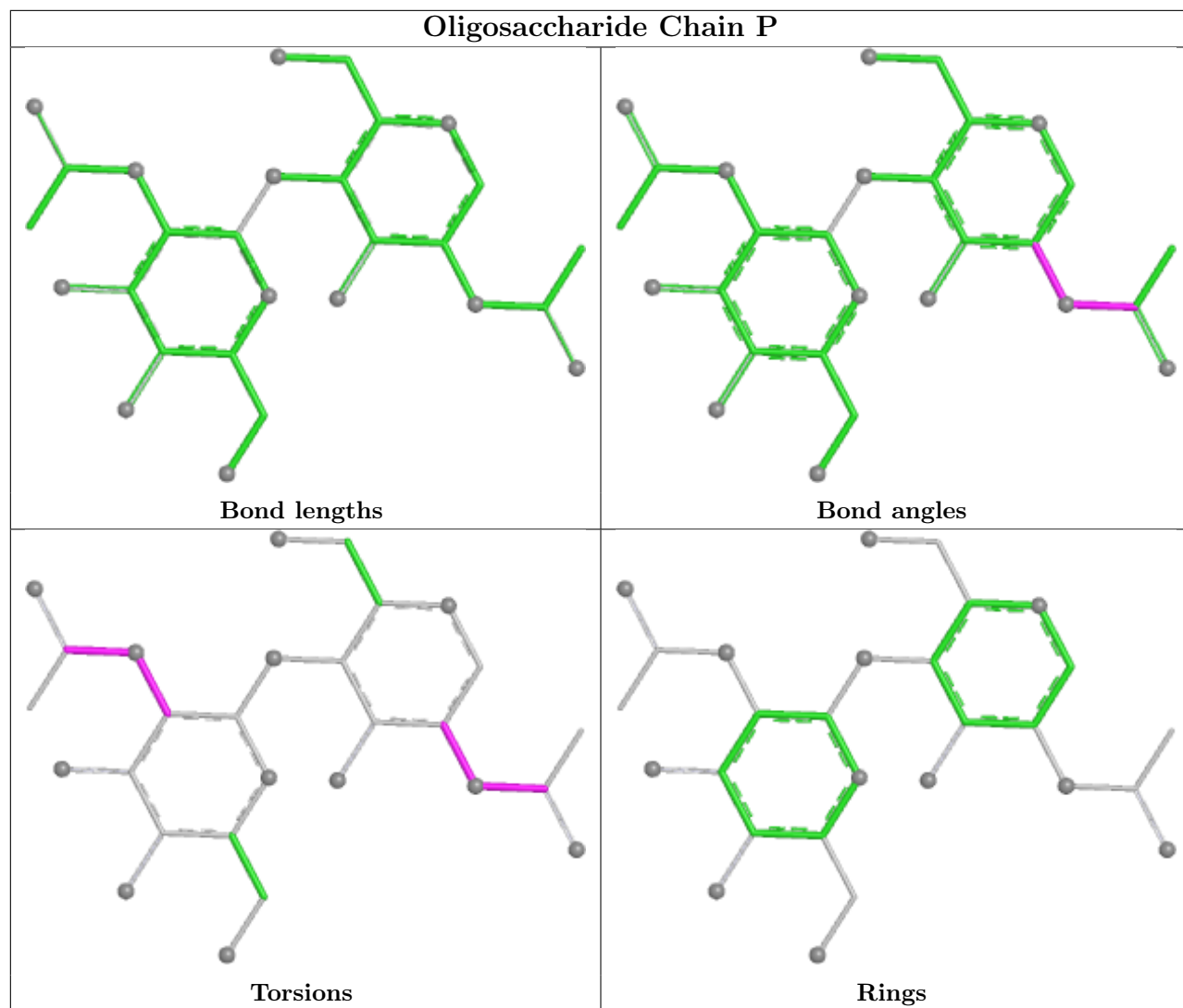


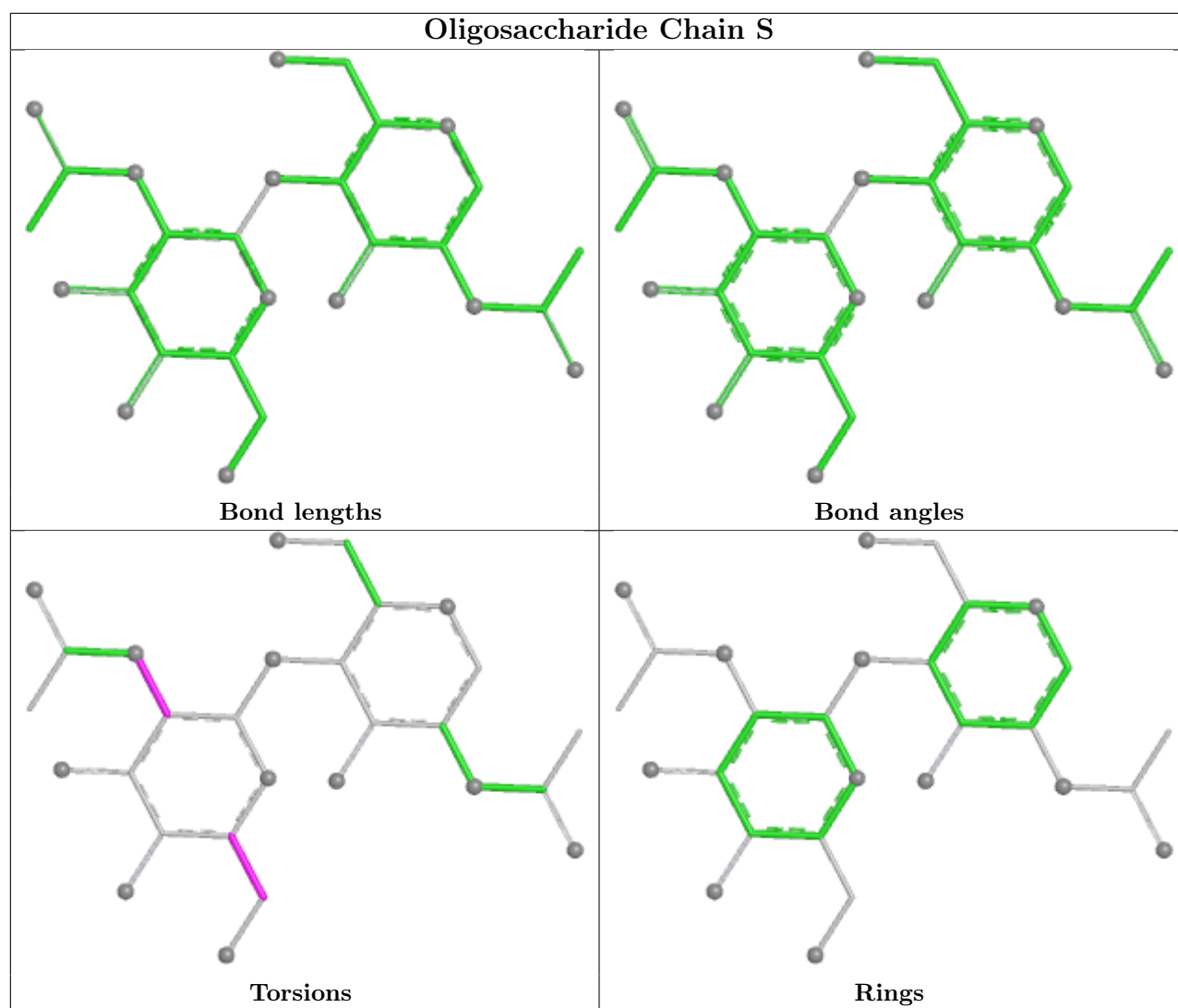


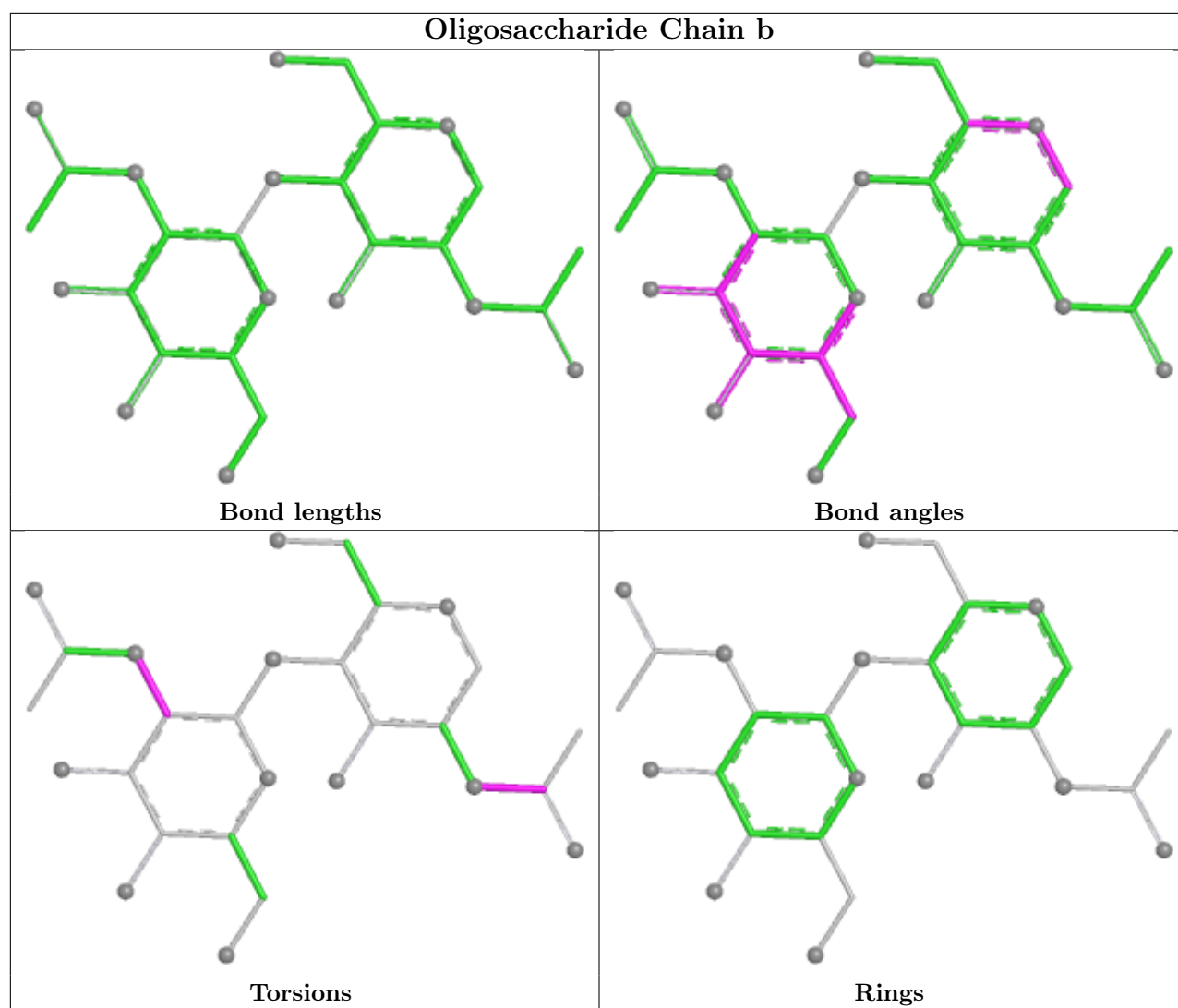


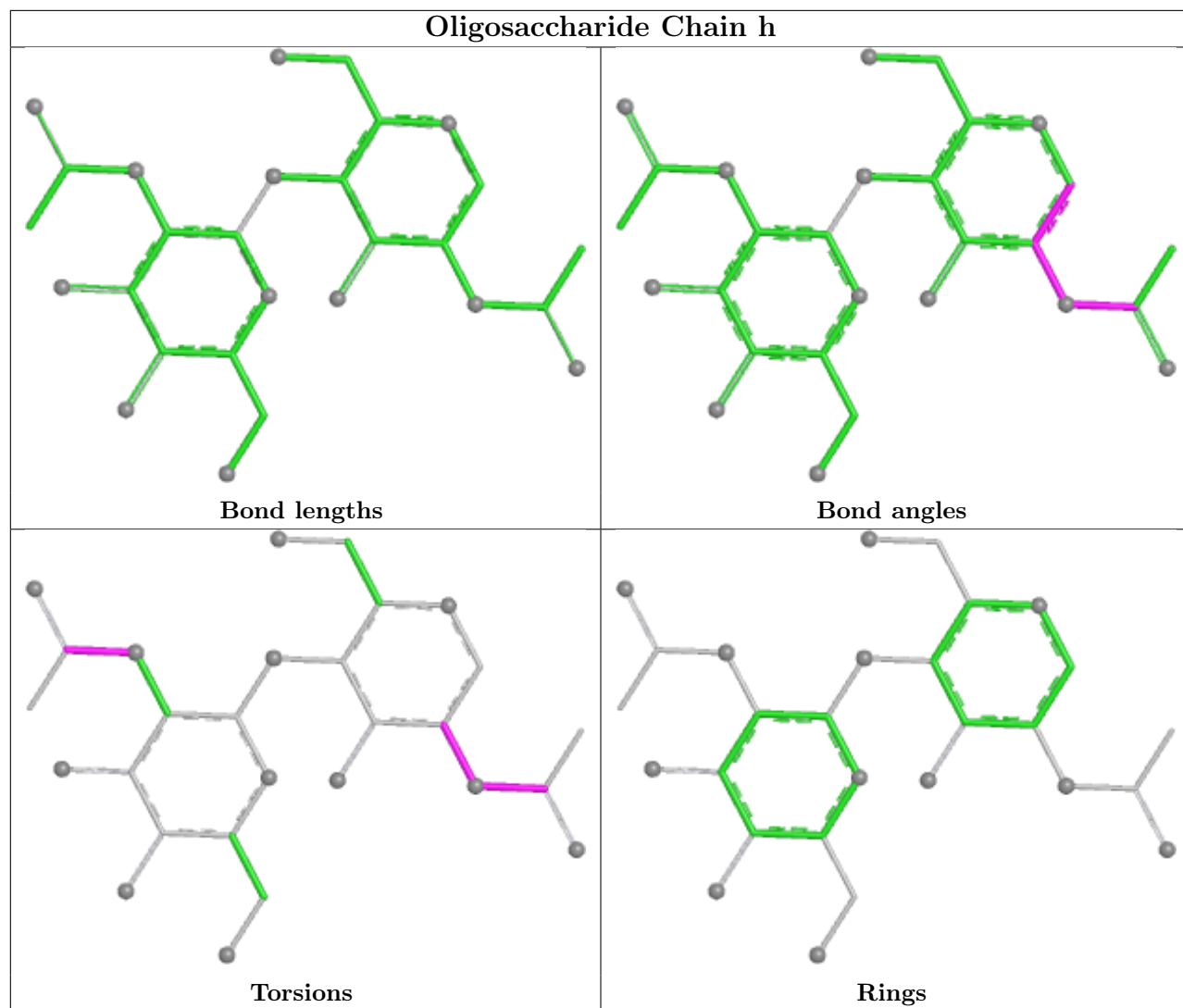


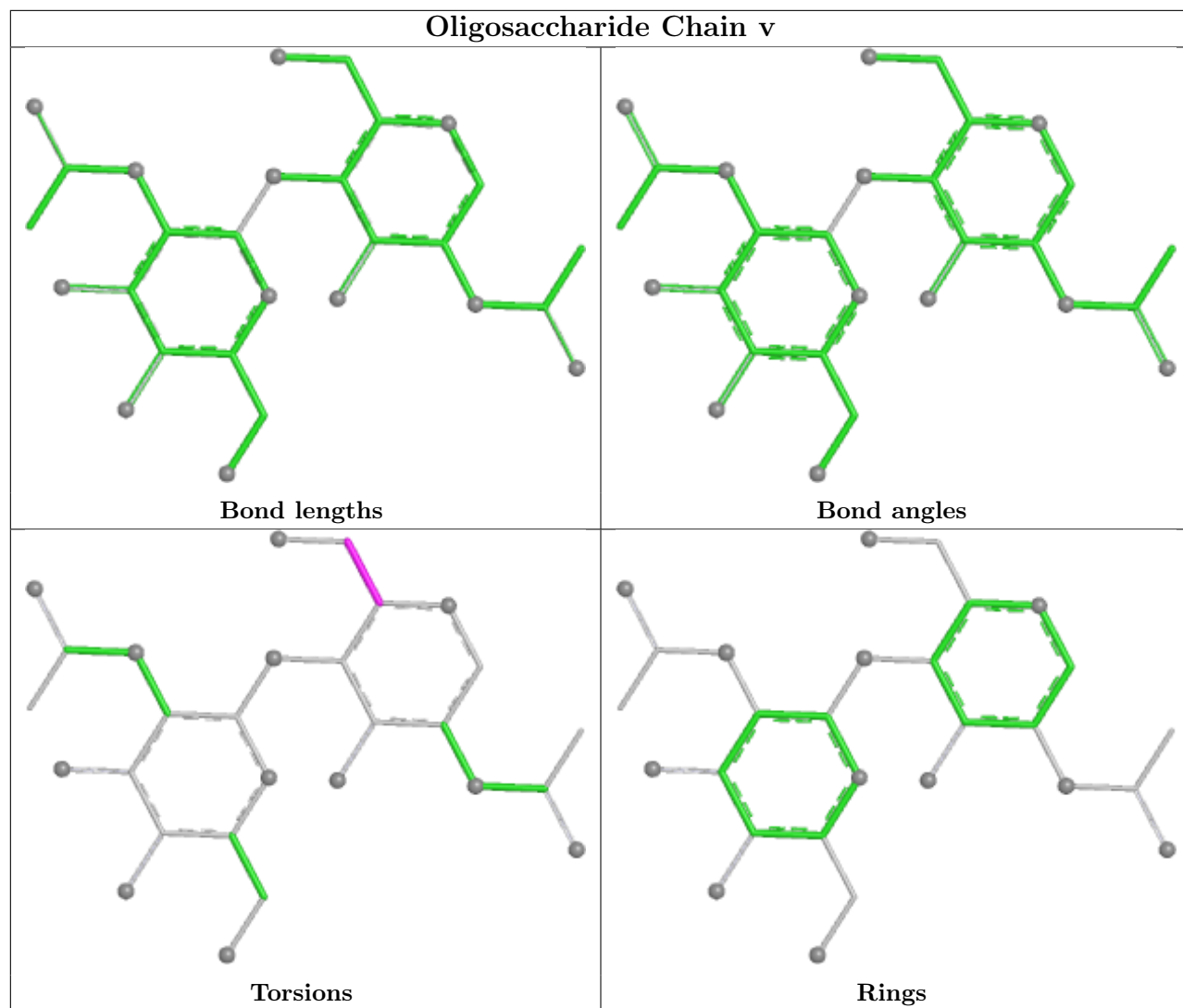


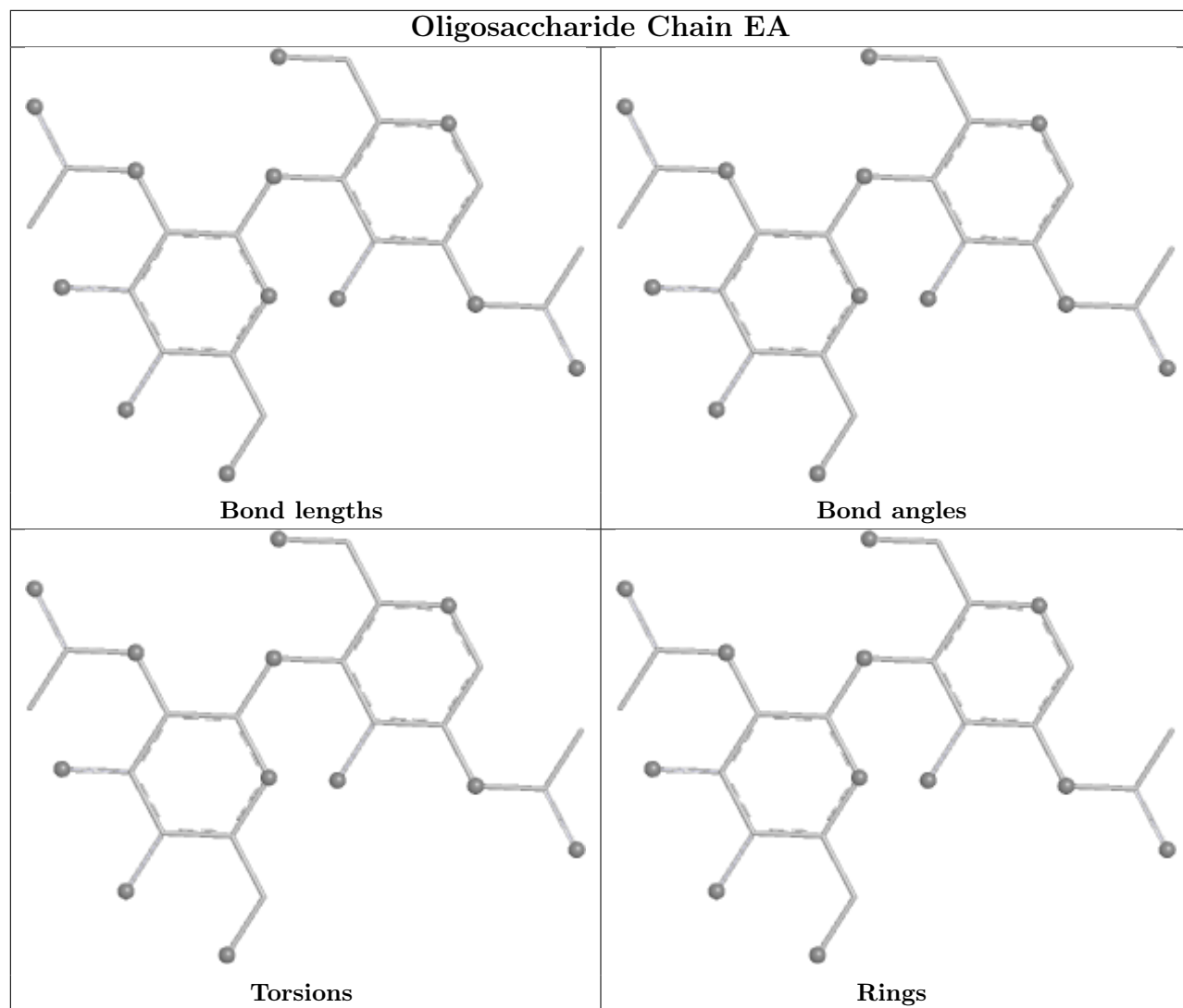


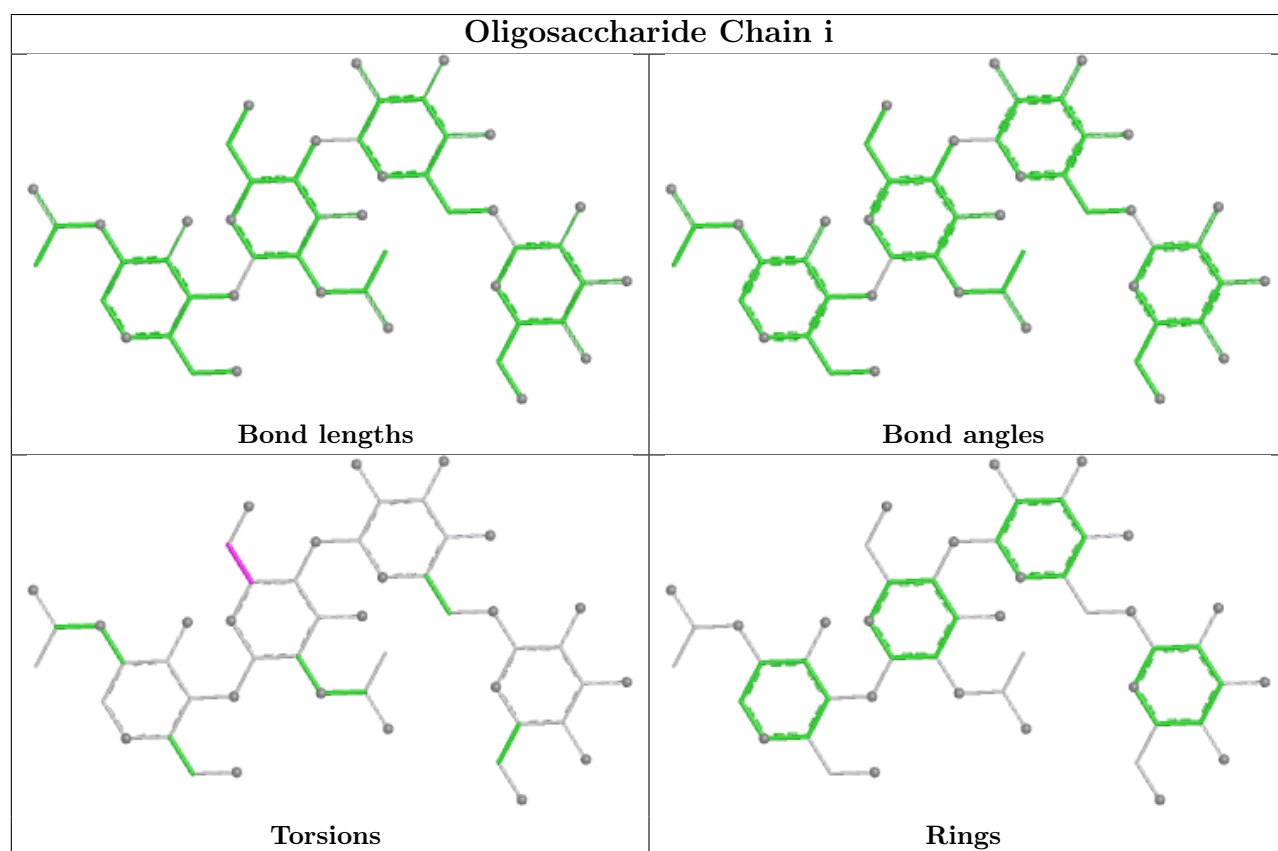
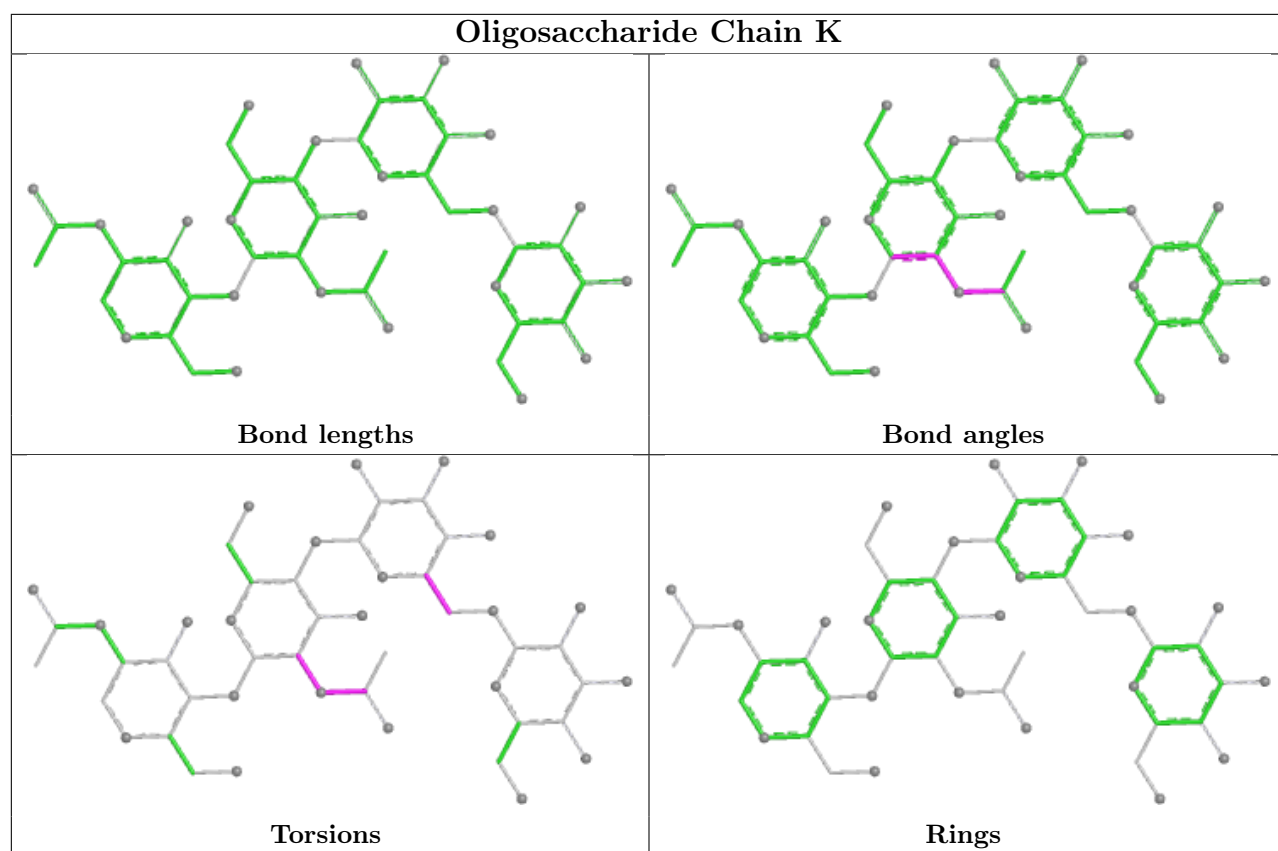


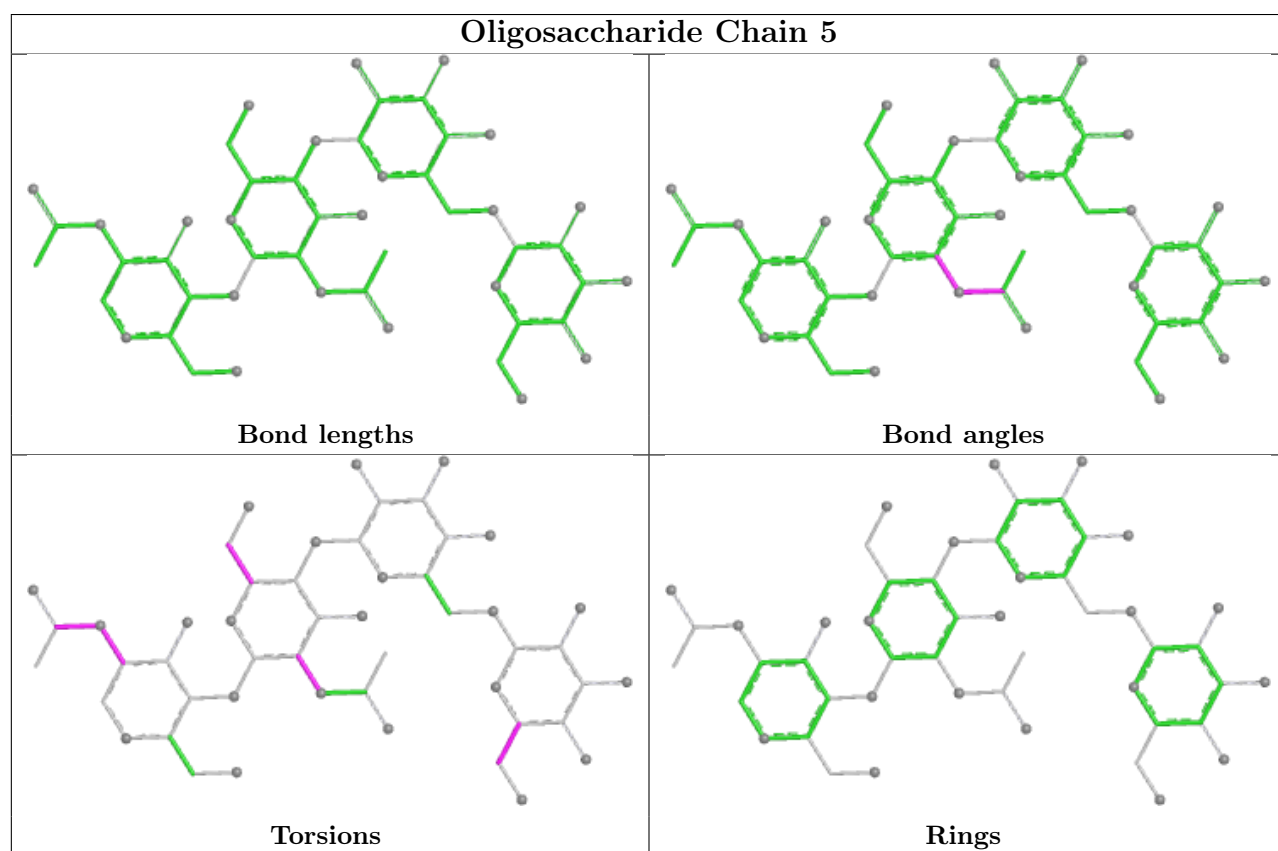
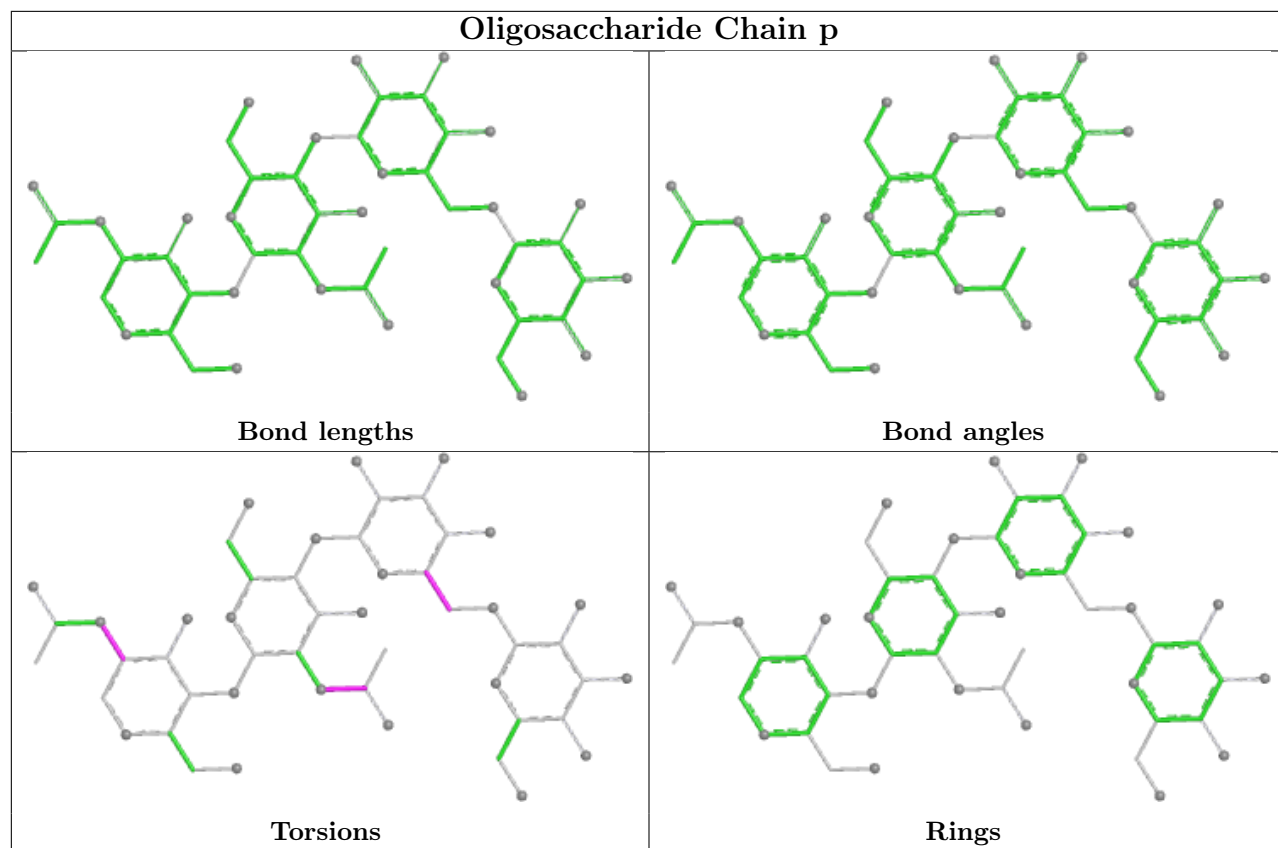


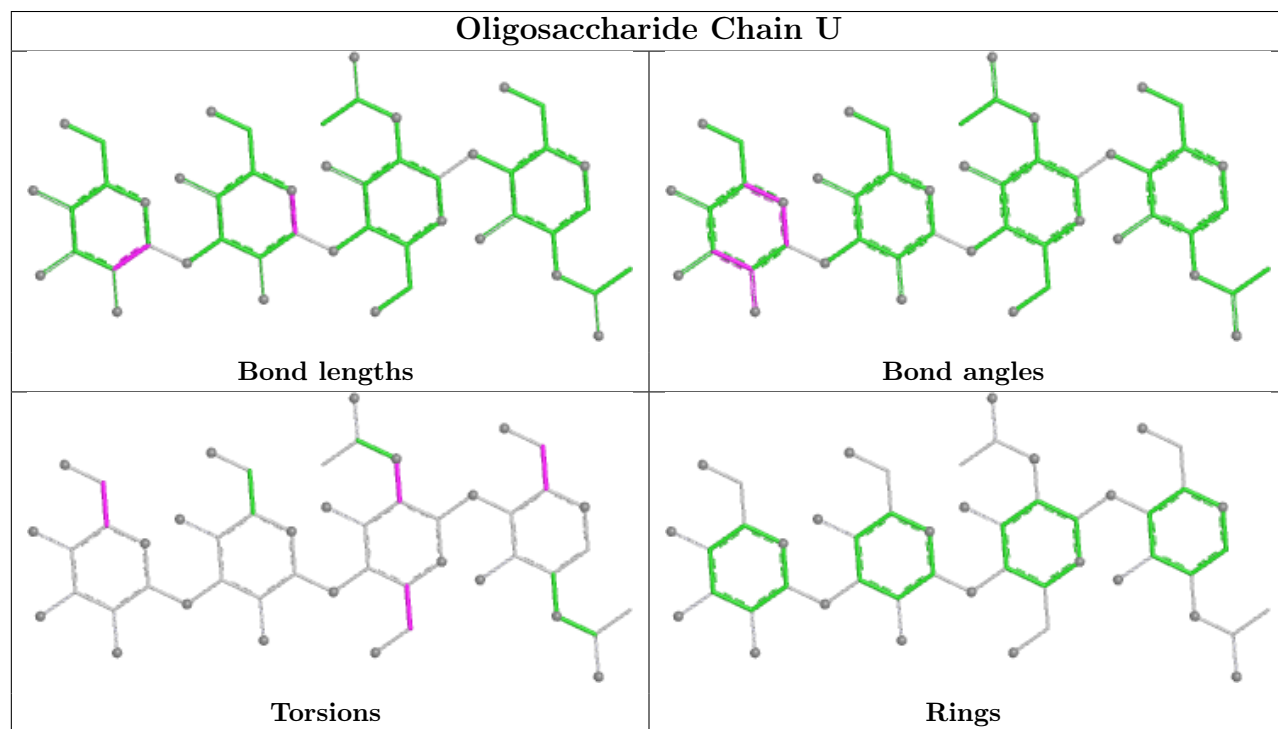
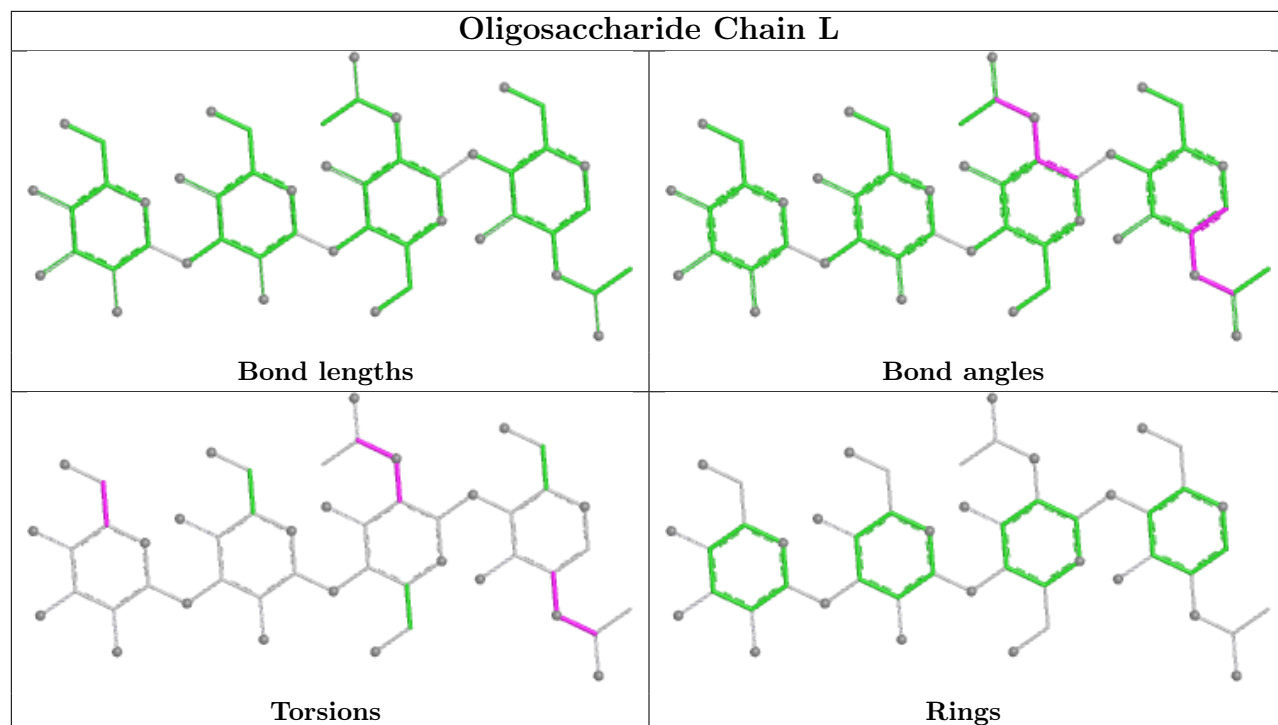


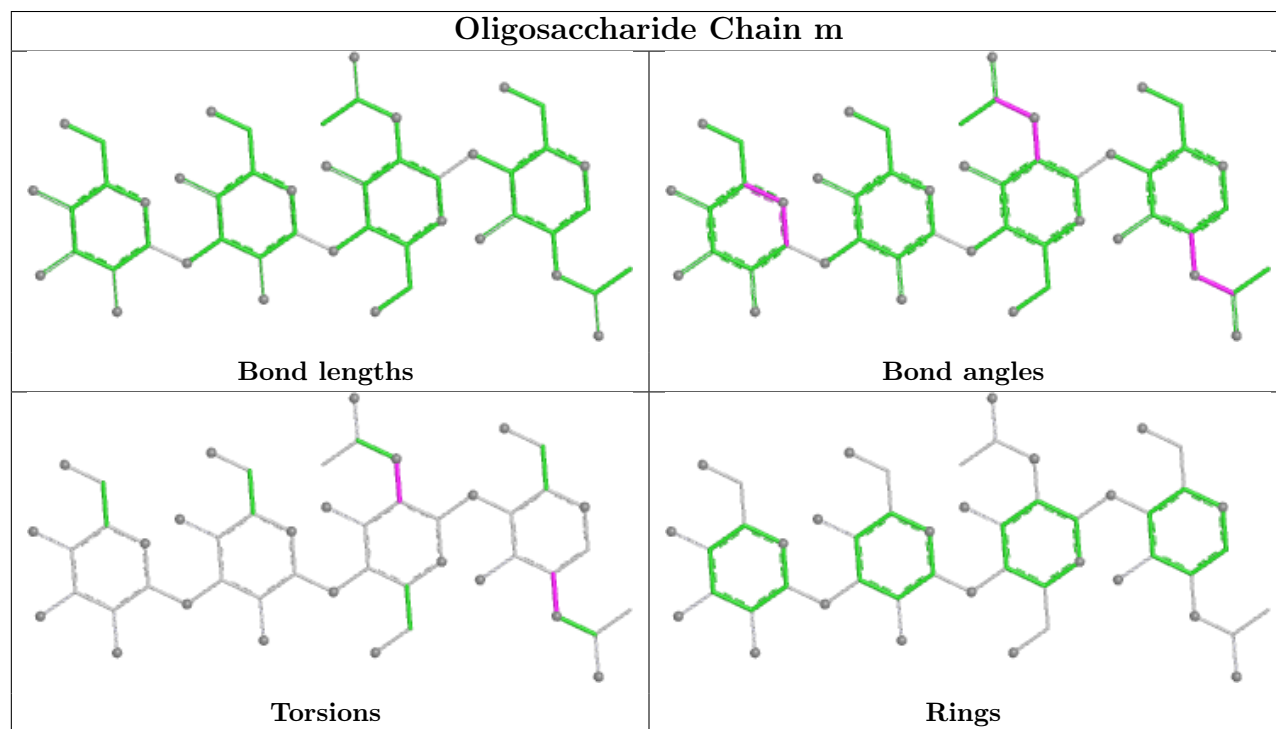
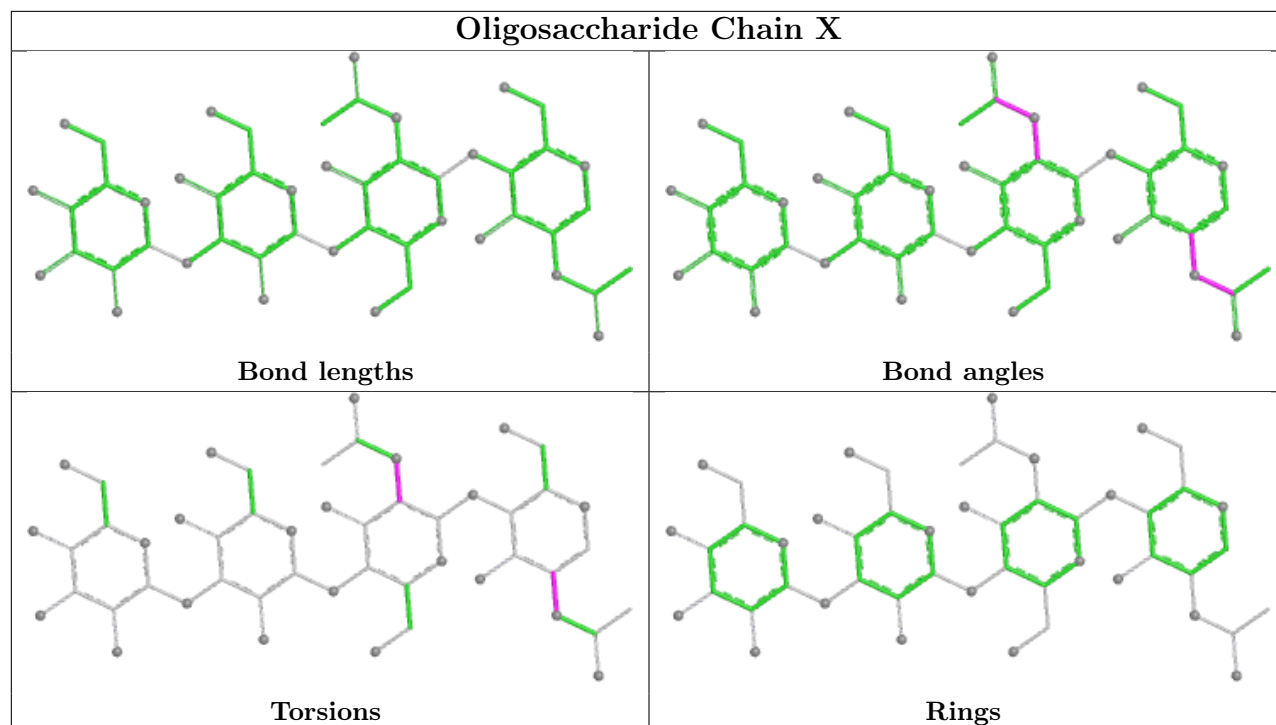


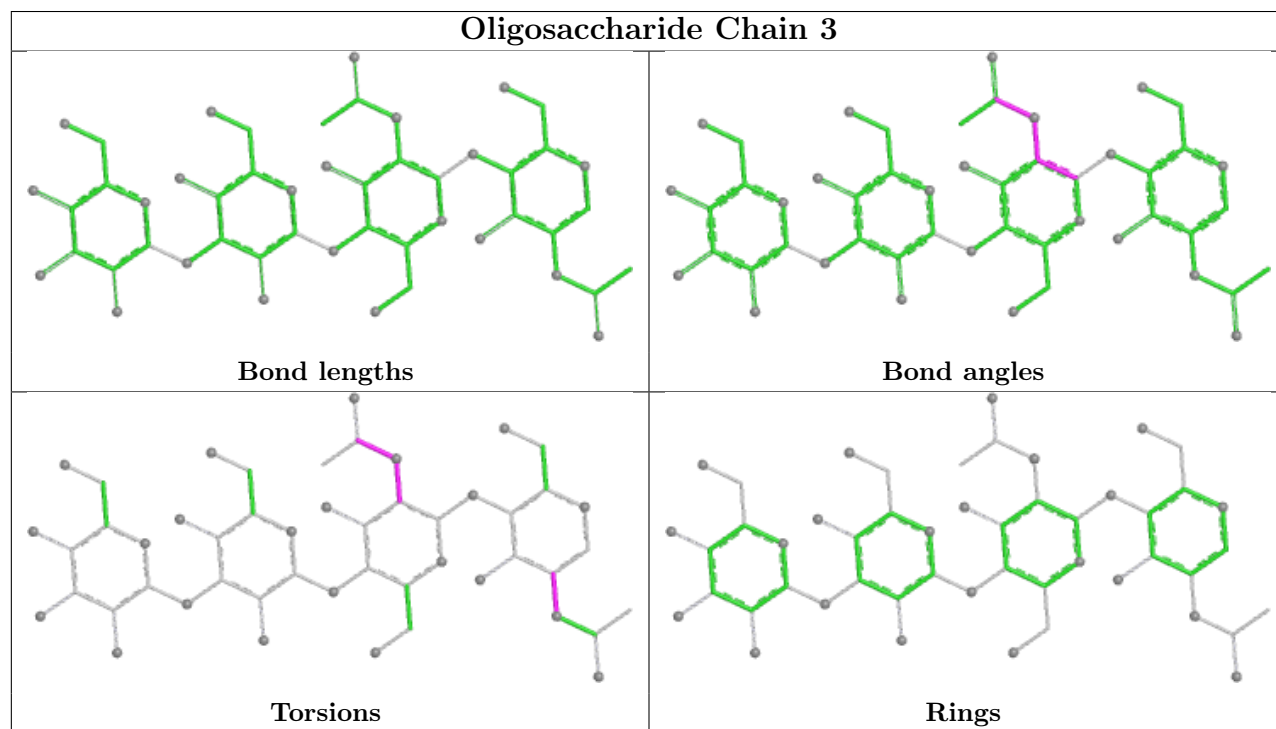
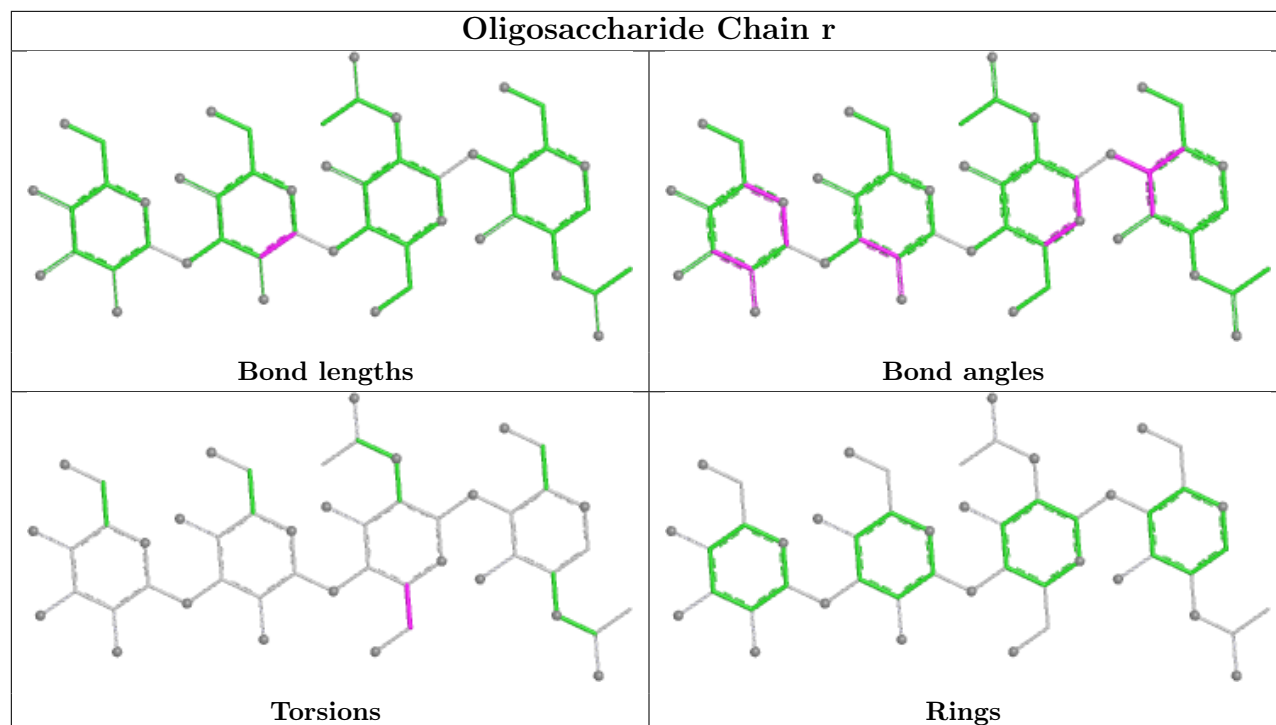




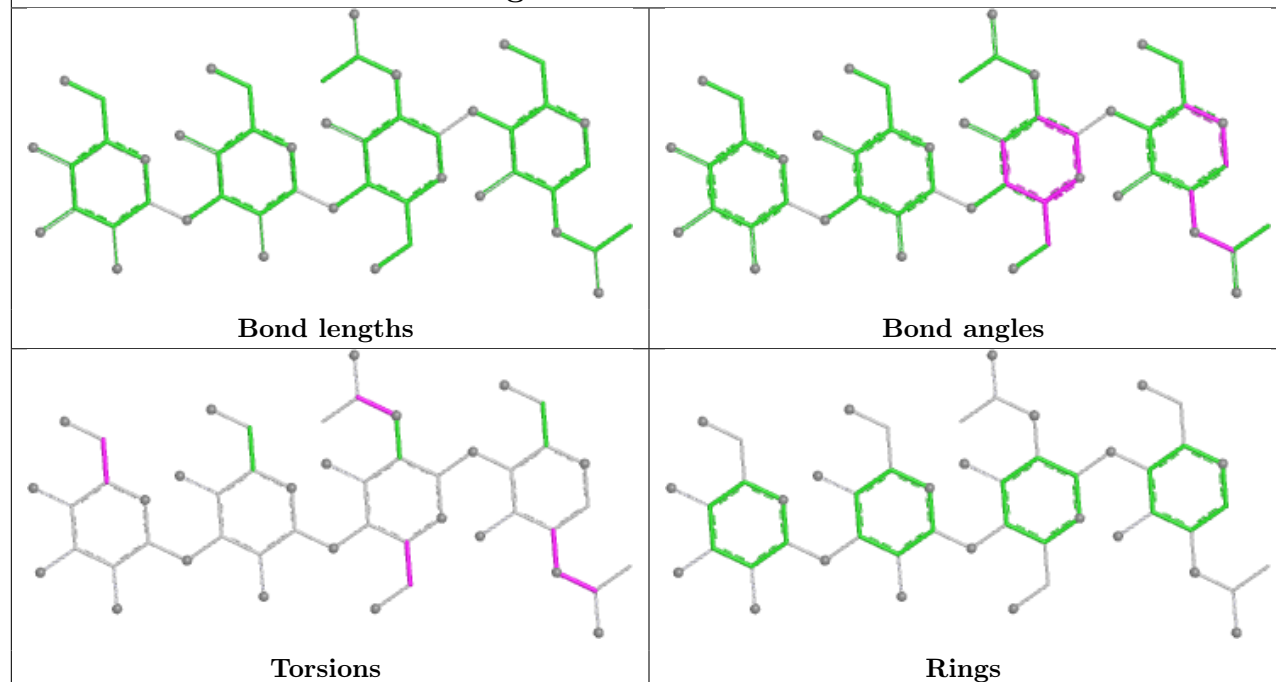




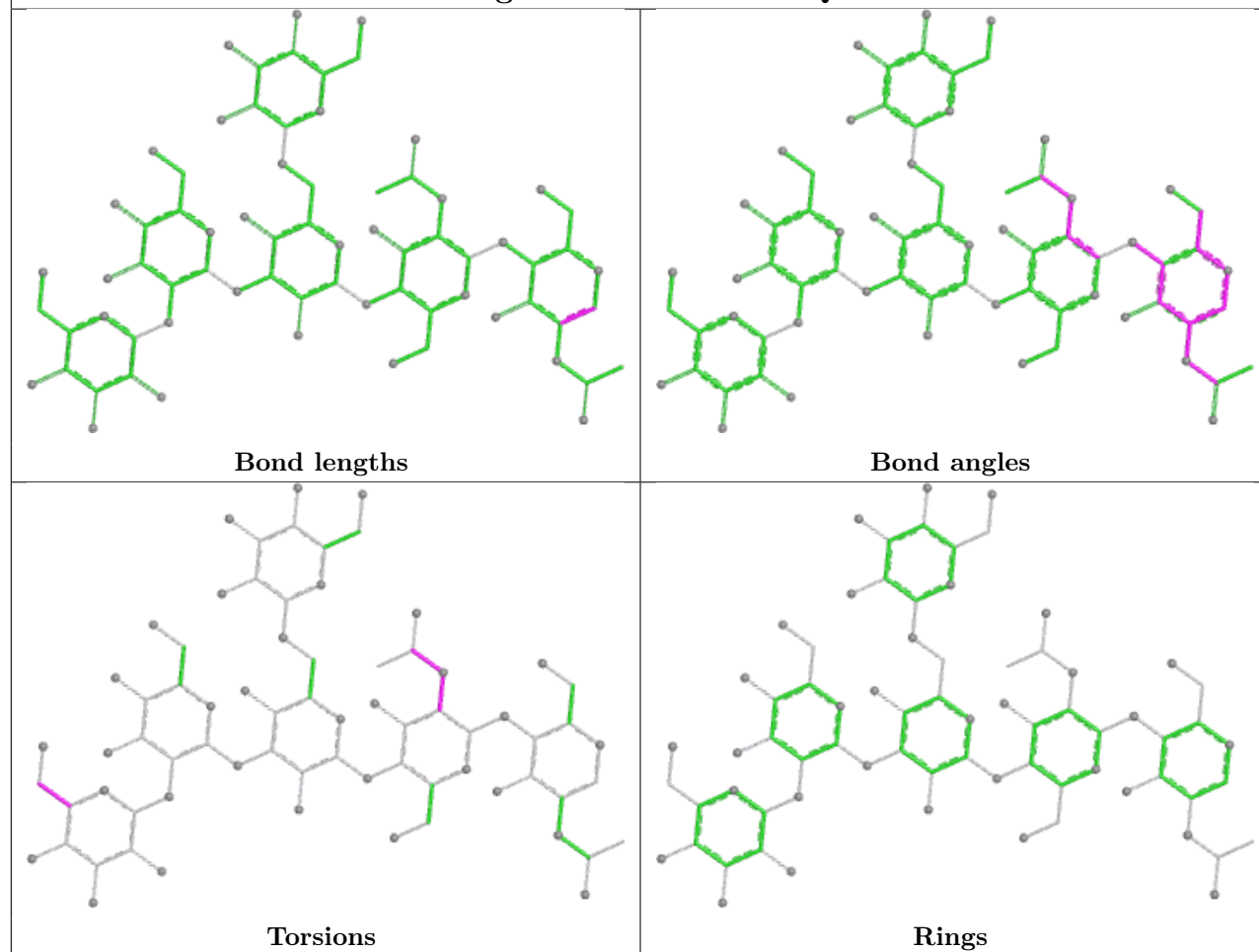


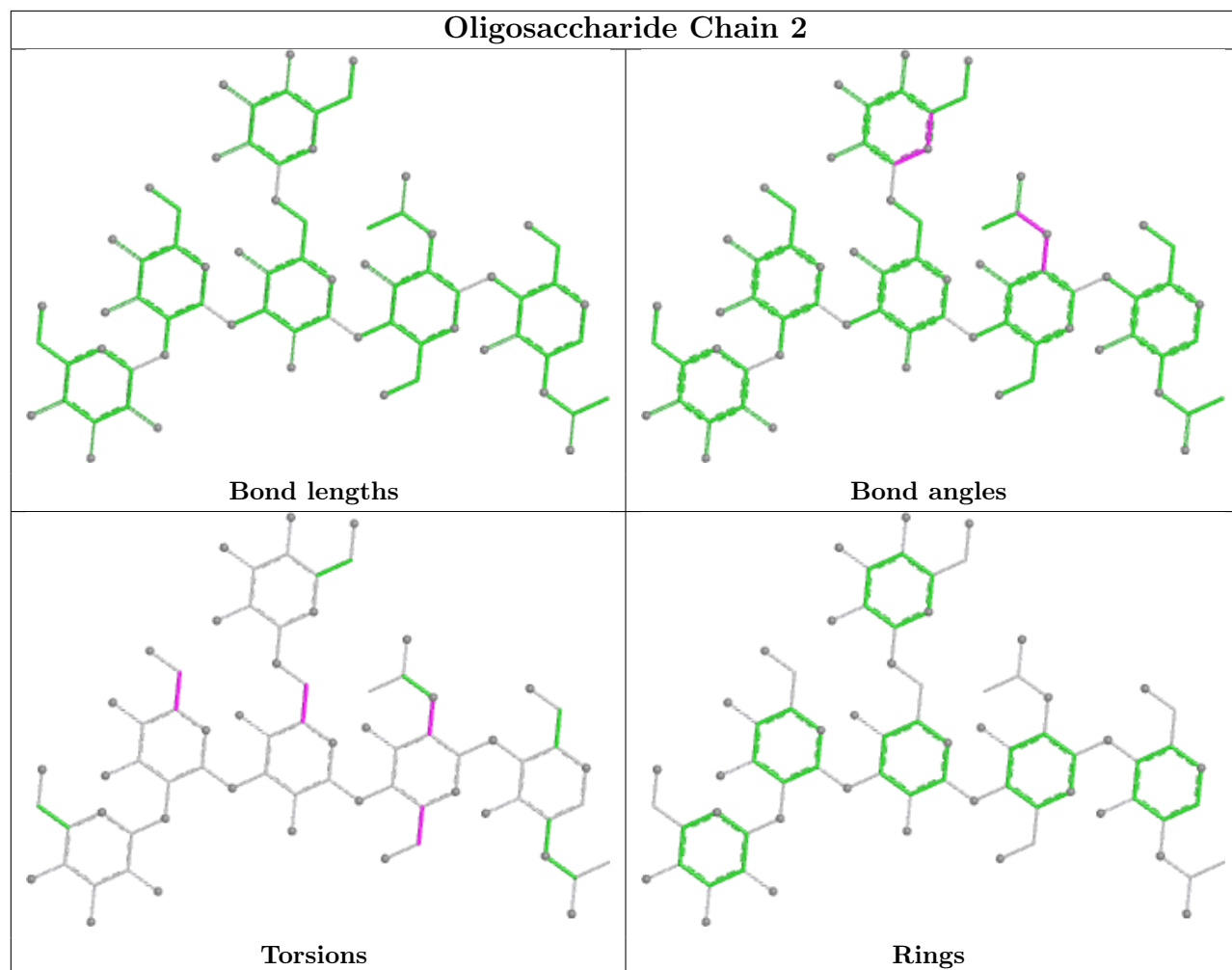


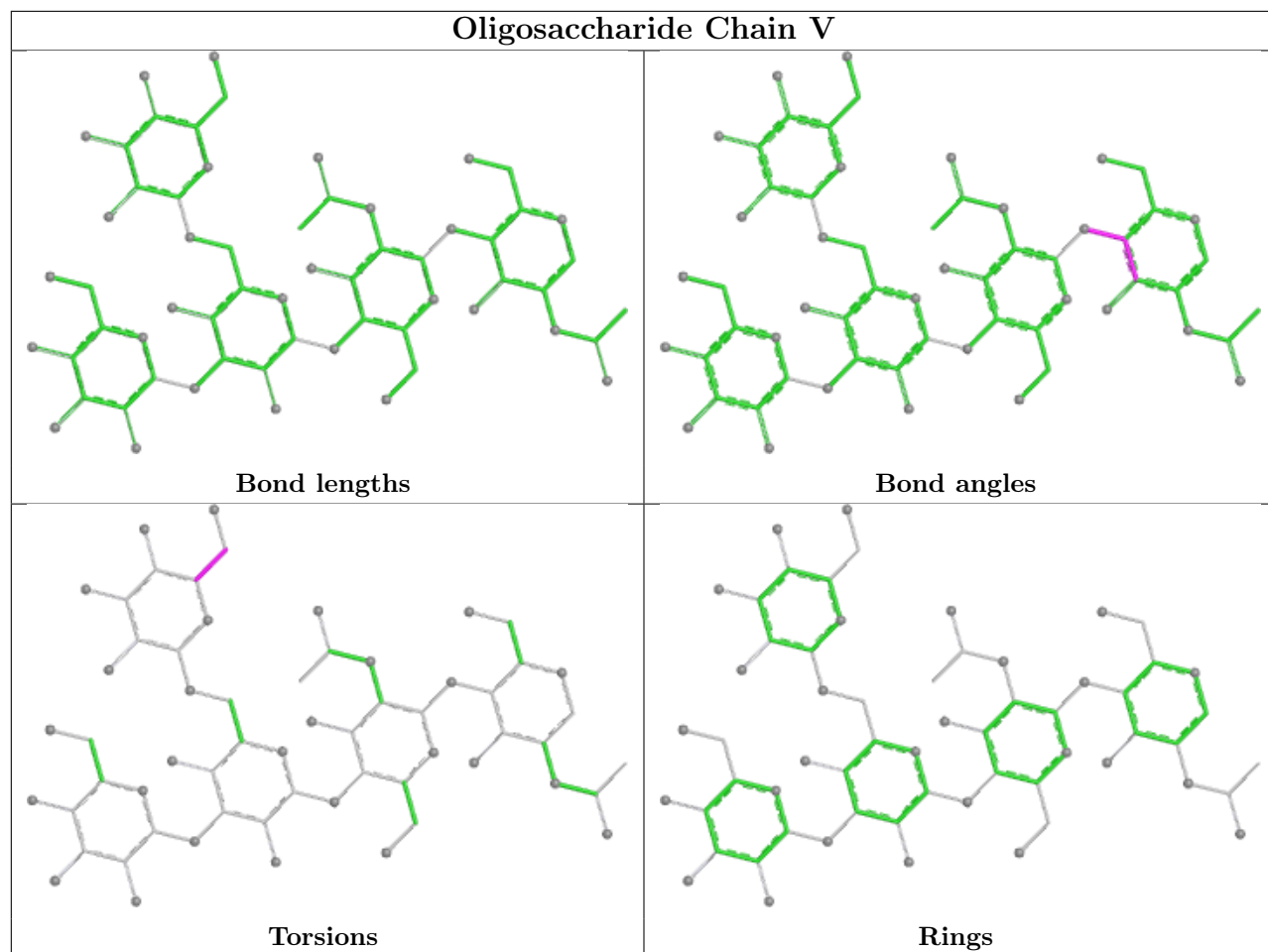
Oligosaccharide Chain 9

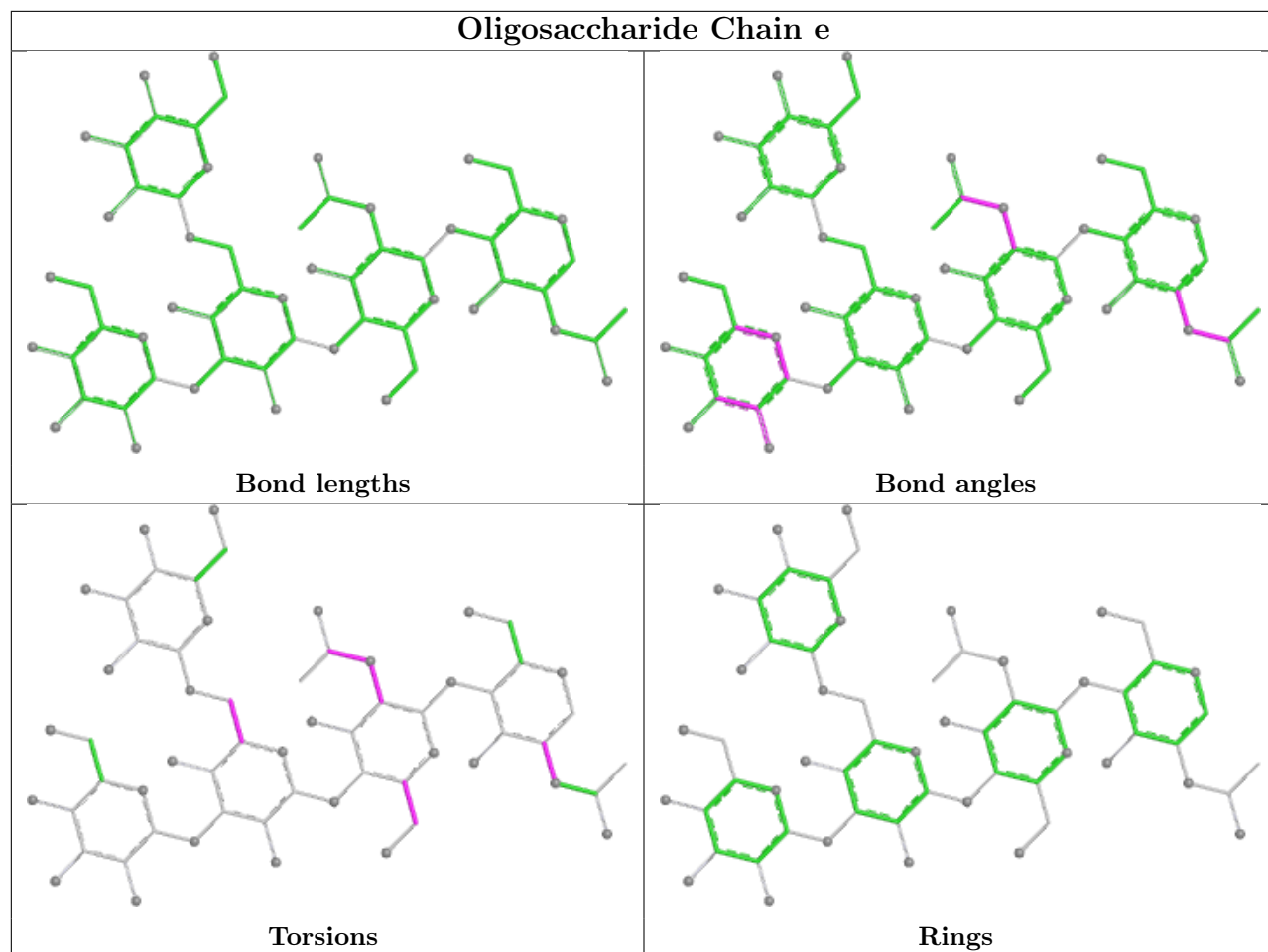


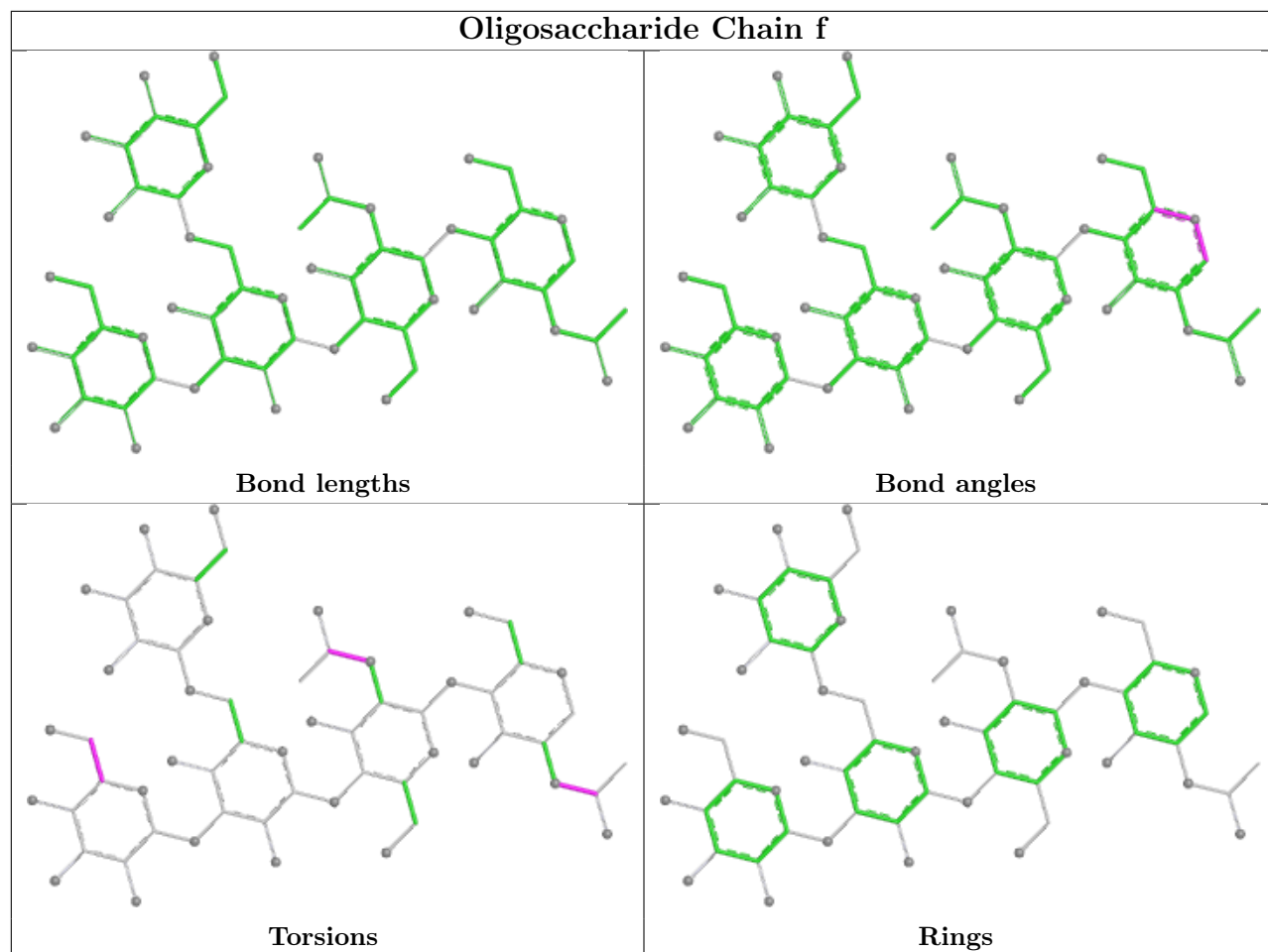
Oligosaccharide Chain Q

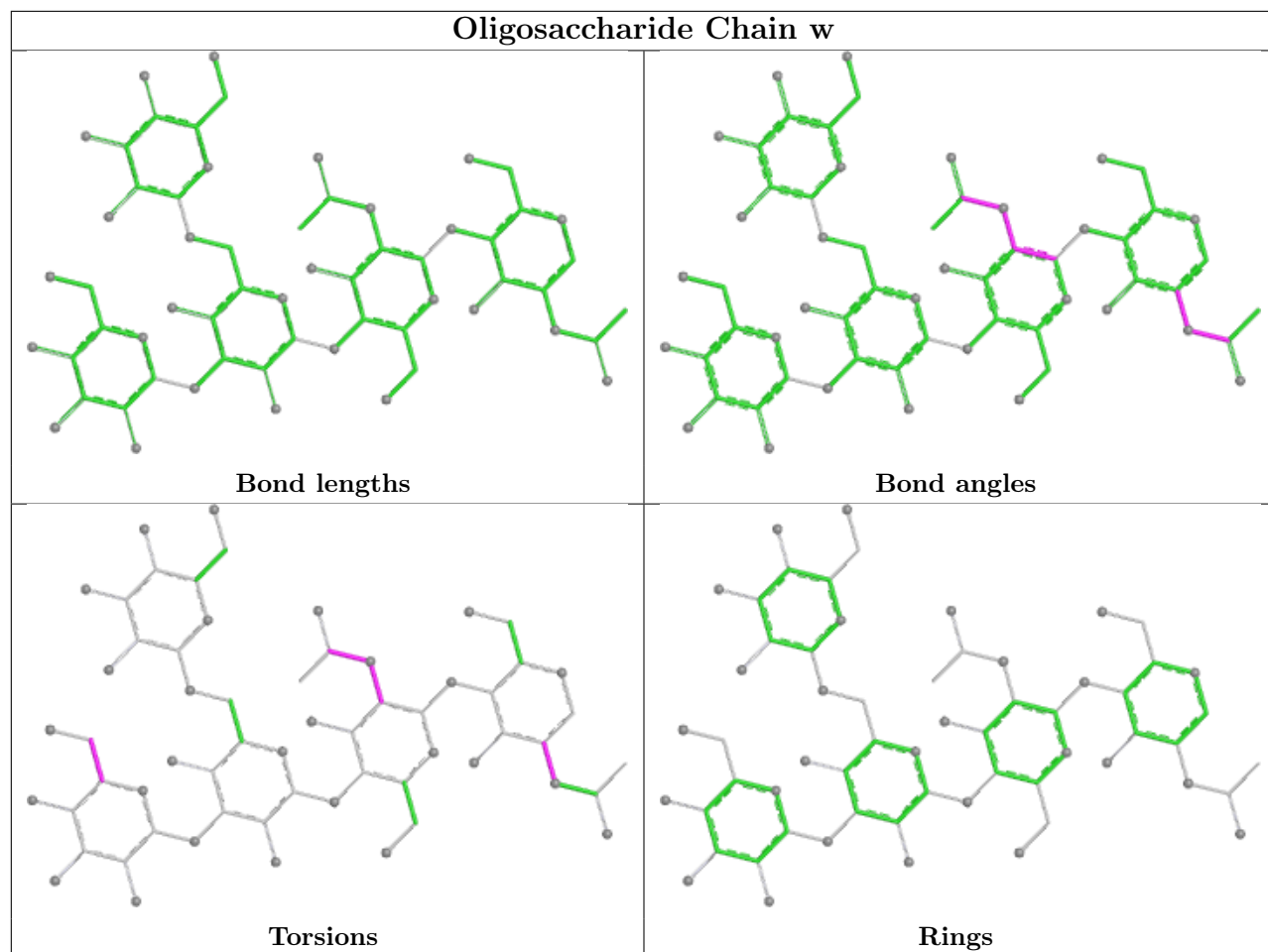


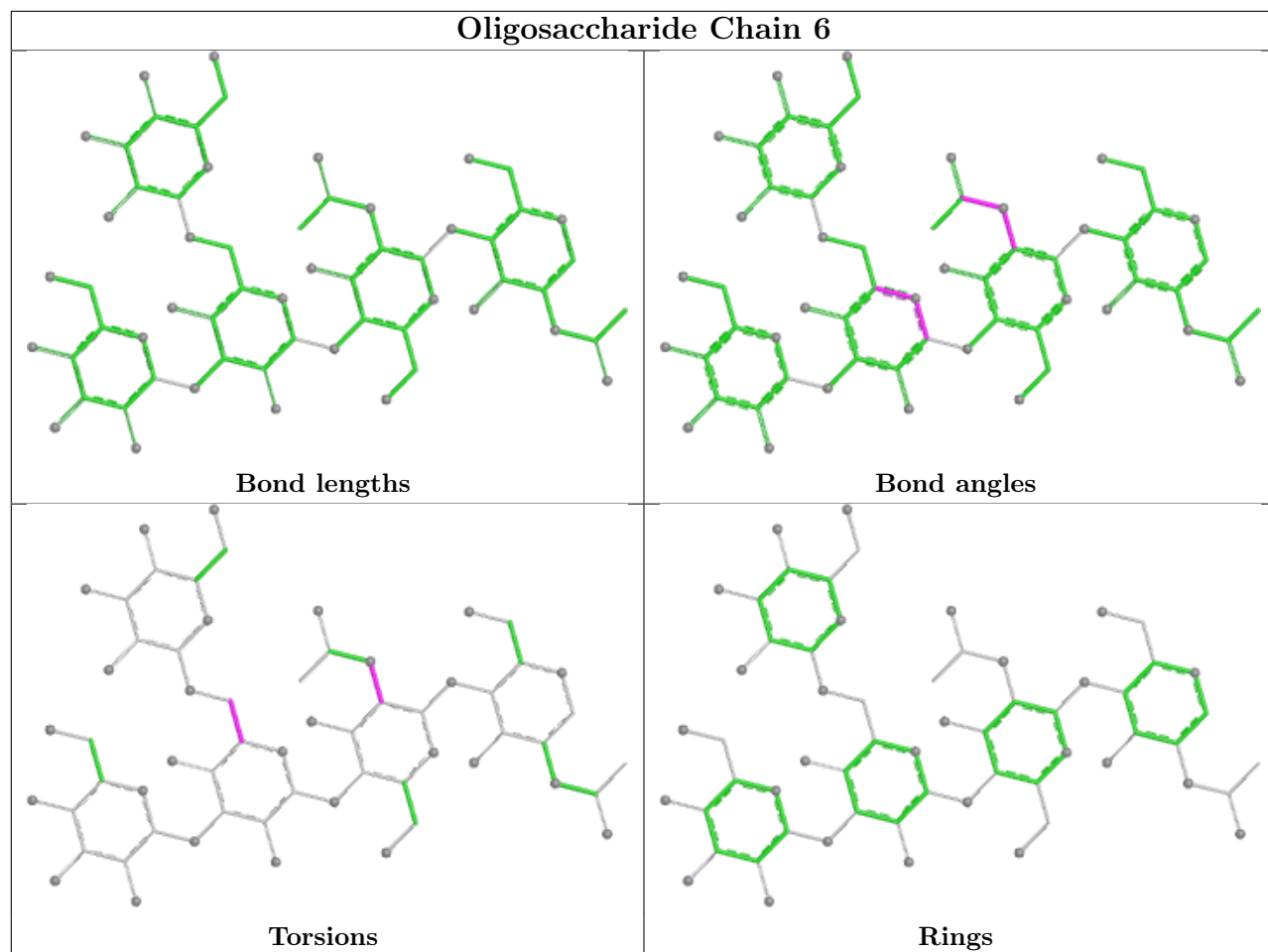


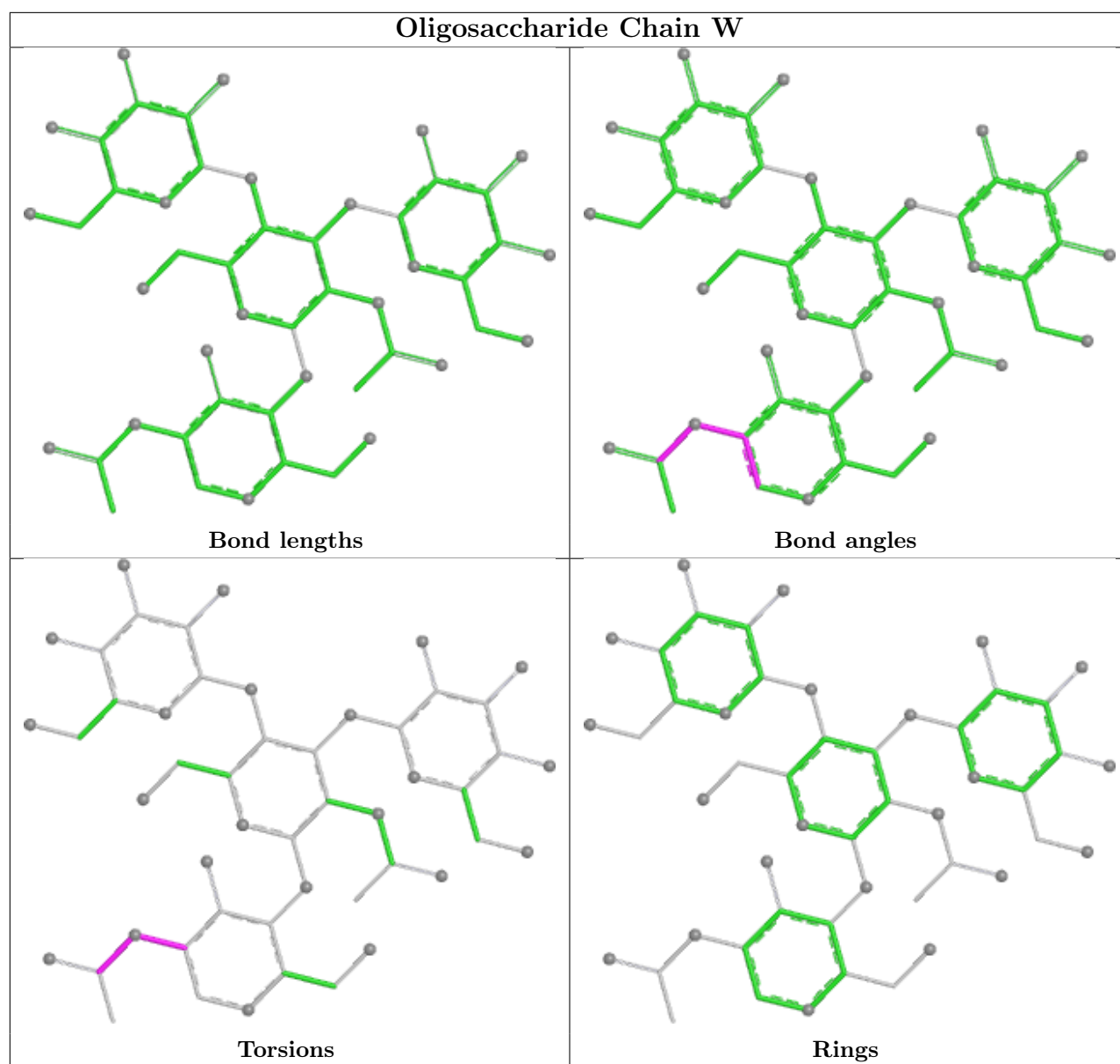


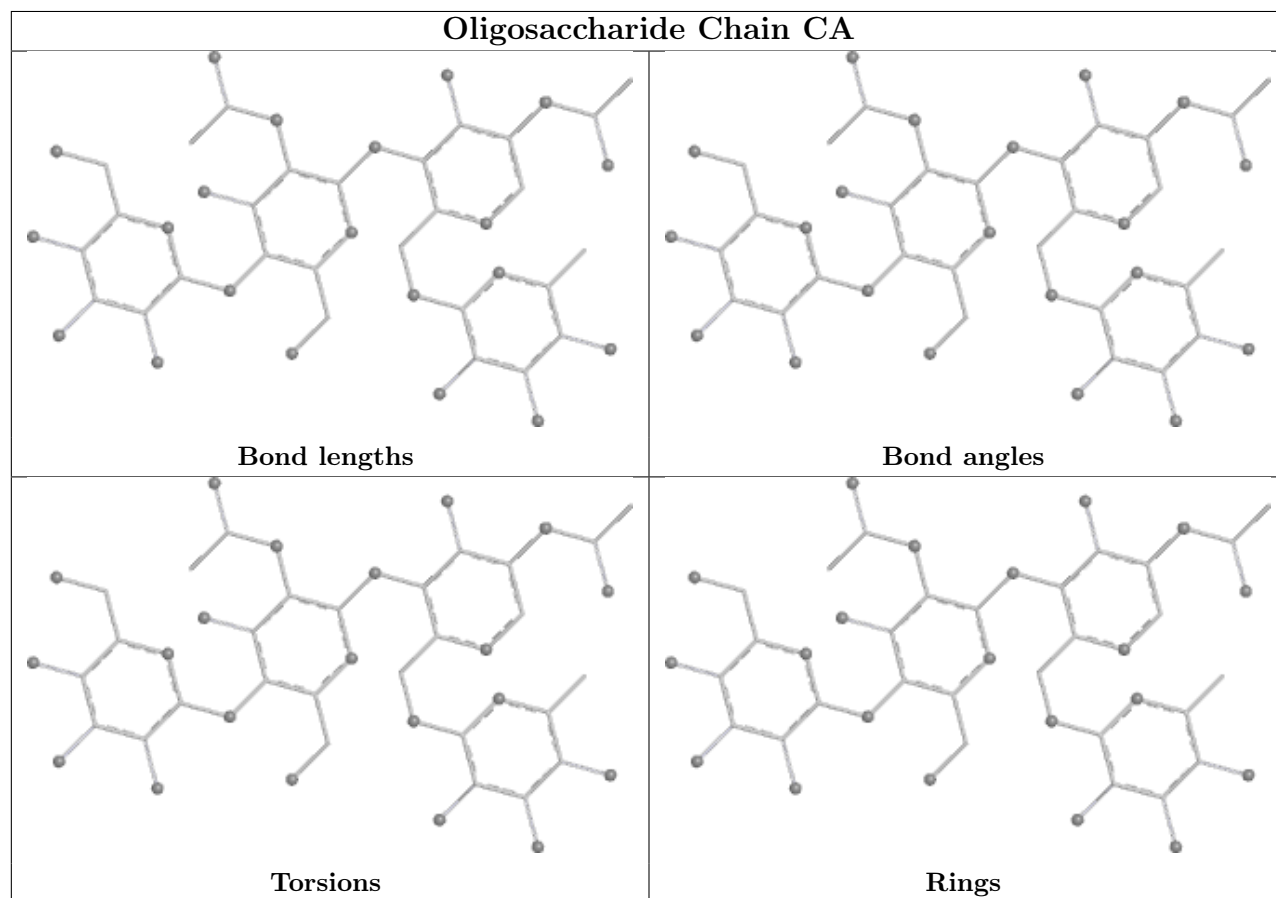
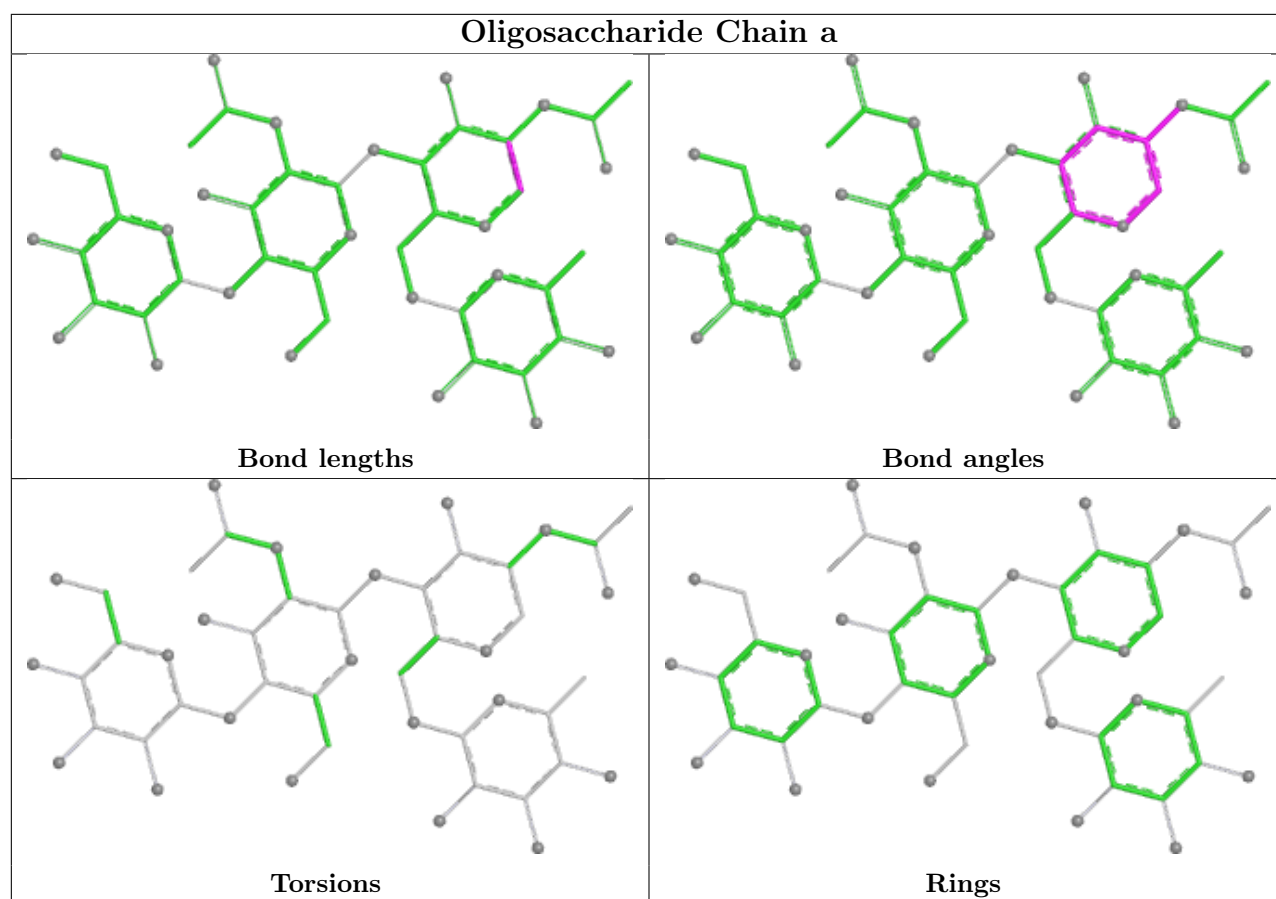


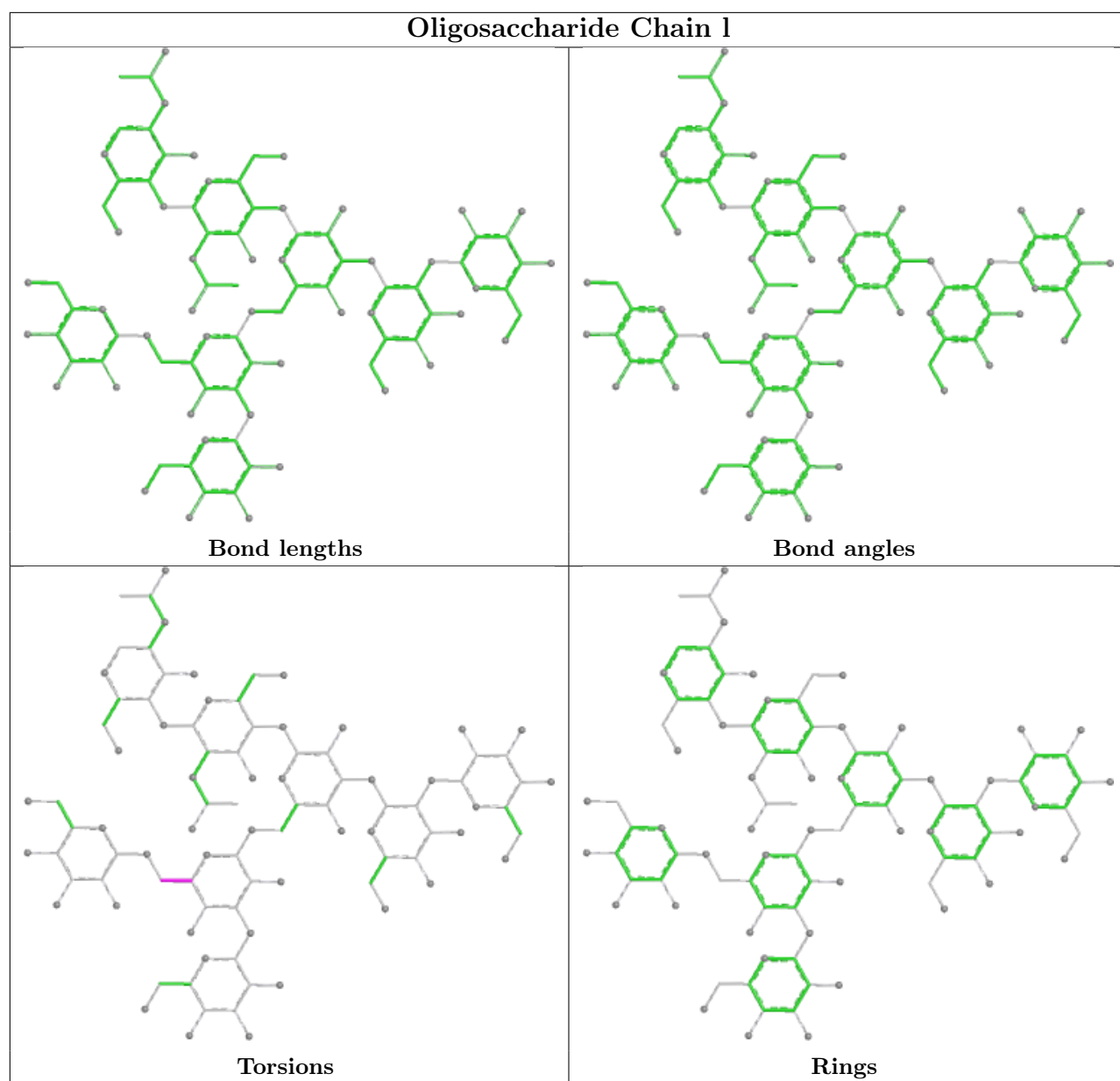


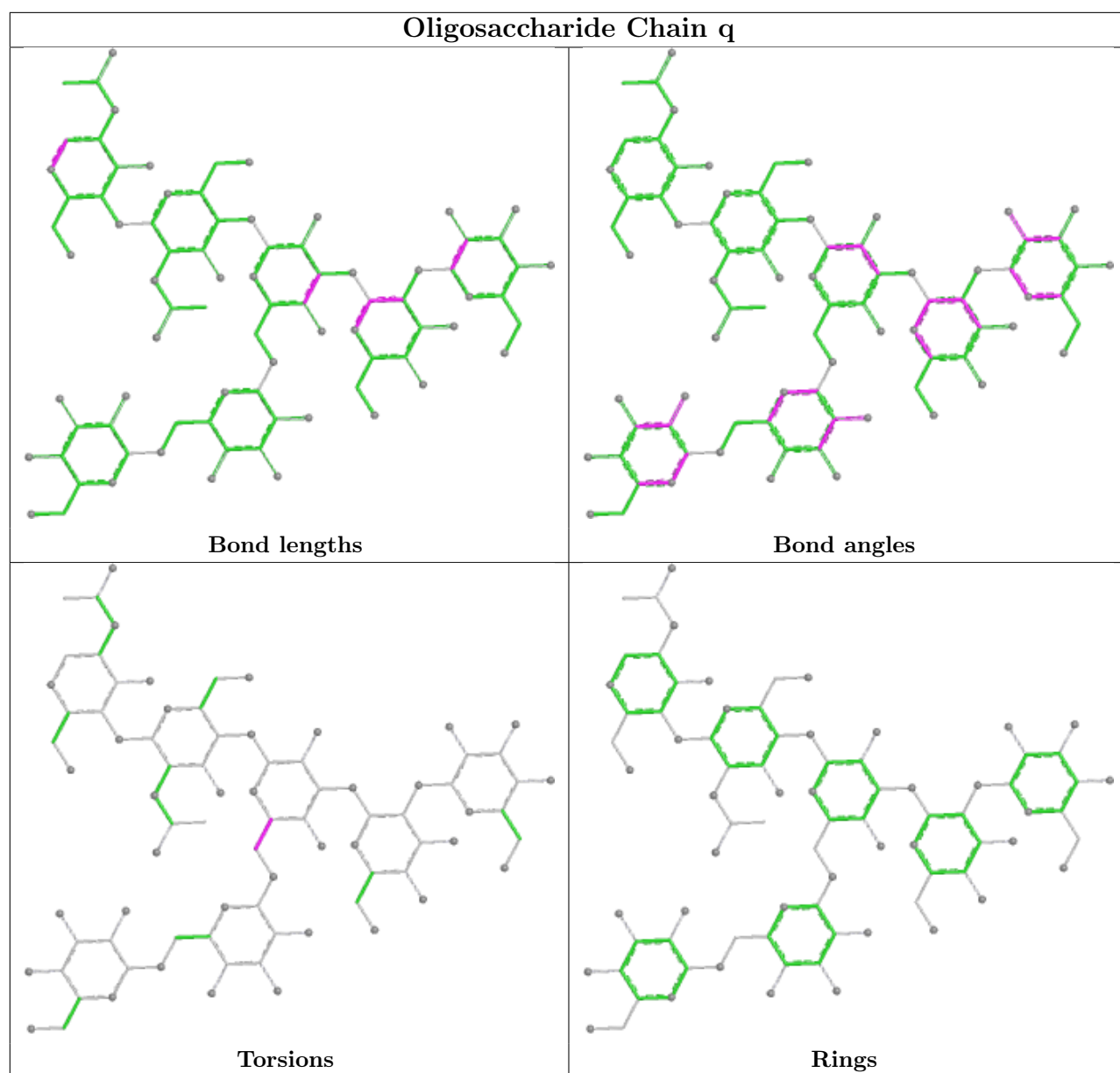


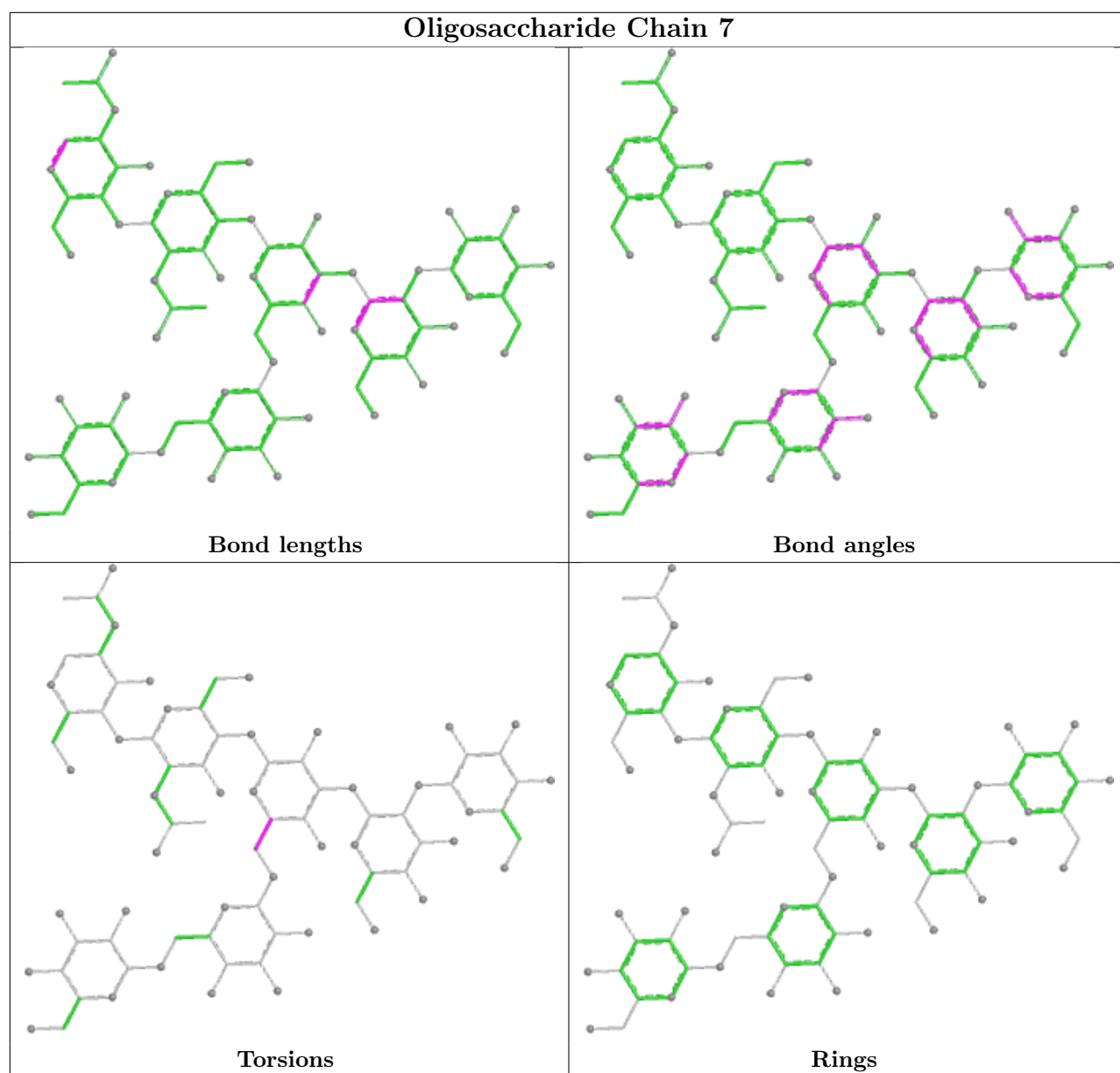


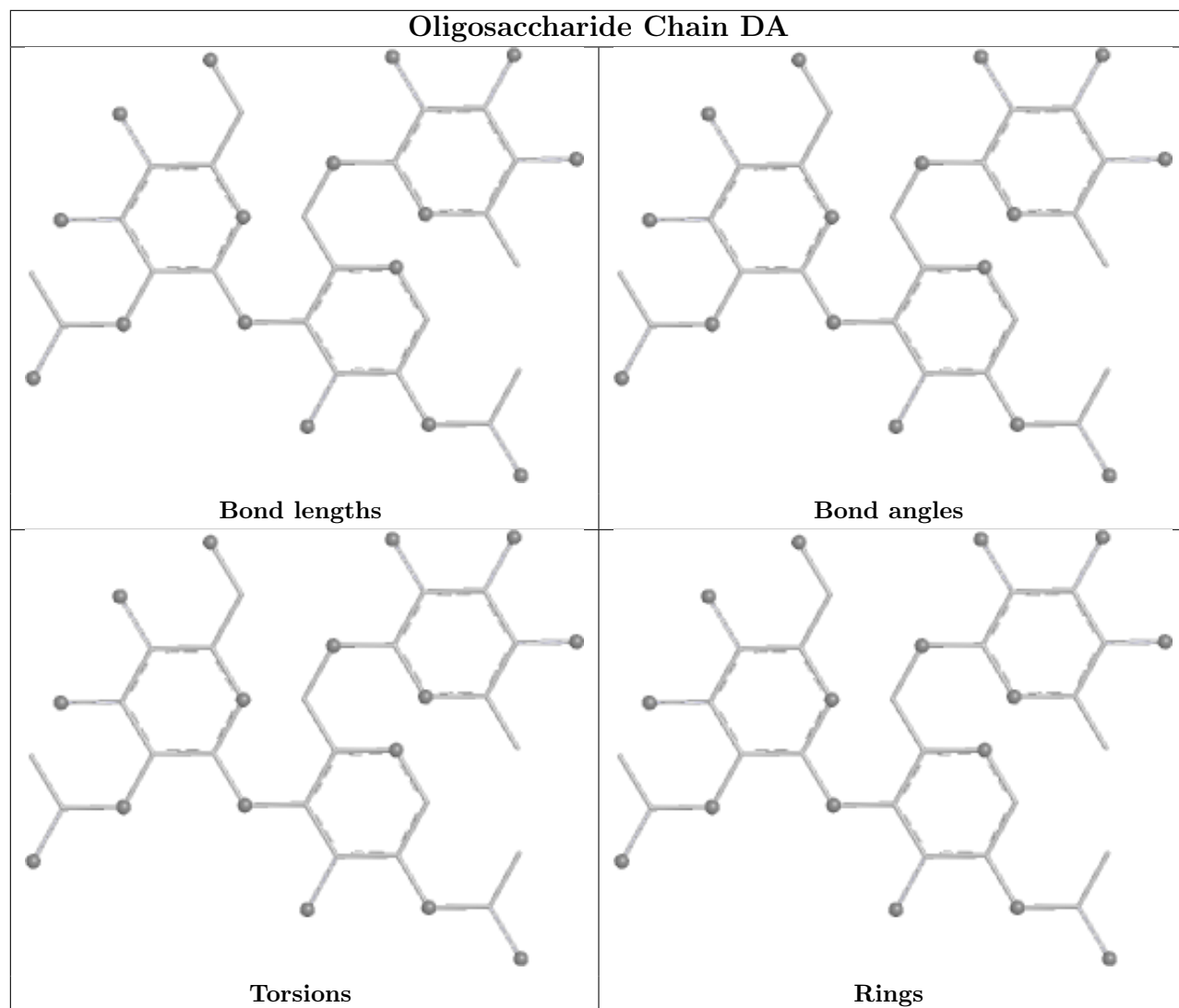


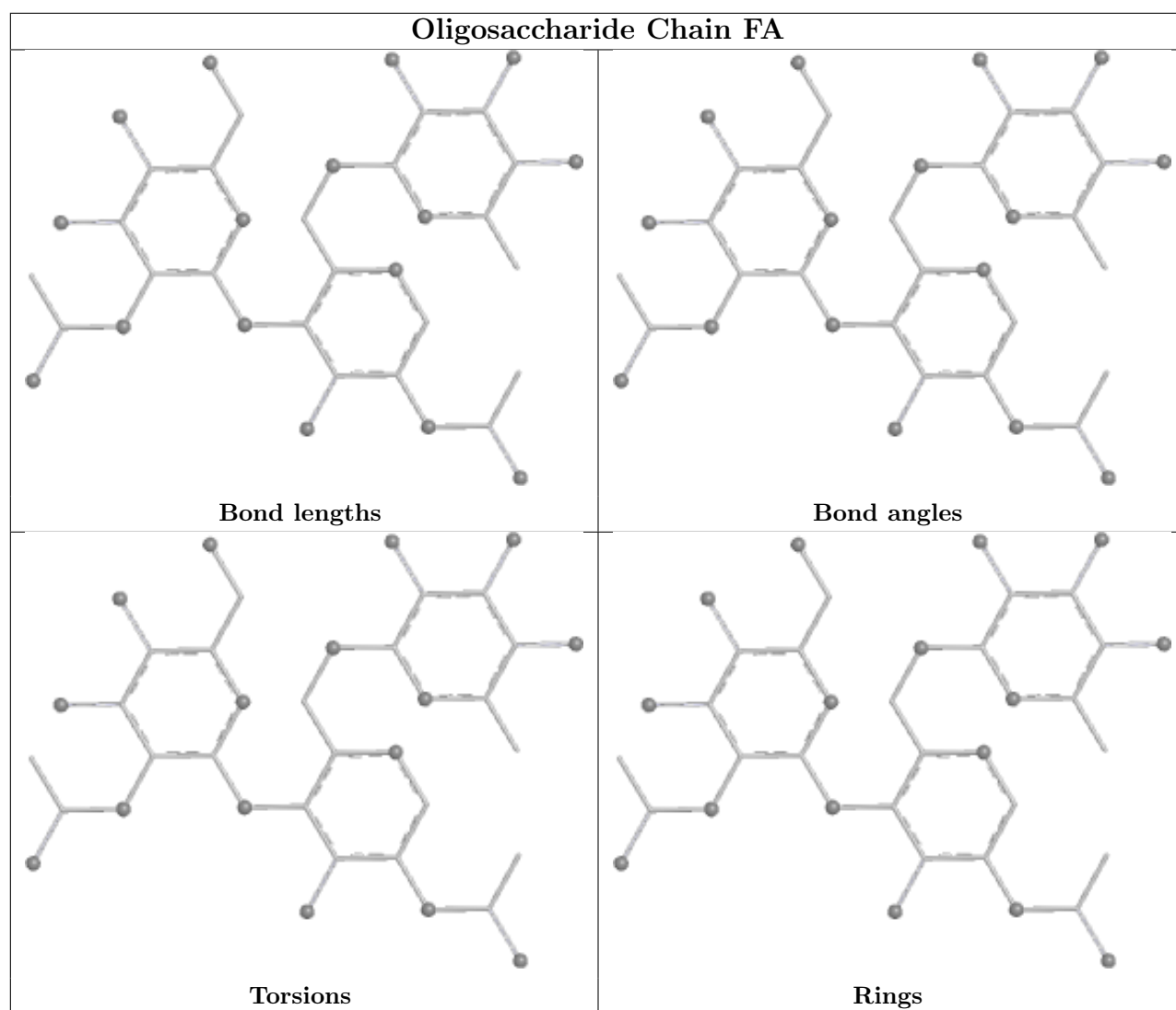












5.6 Ligand geometry [i](#)

14 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	$\# Z > 2$
13	NAG	A	803	1	14,14,15	0.35	0	17,19,21	0.51	0
13	NAG	C	802	1	14,14,15	0.38	0	17,19,21	0.44	0
13	NAG	C	805	1	14,14,15	0.36	0	17,19,21	1.15	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
14	83G	E	801	-	33,33,33	0.45	0	44,47,47	1.02	3 (6%)
14	83G	A	801	-	33,33,33	0.41	0	44,47,47	0.82	1 (2%)
14	83G	C	801	-	33,33,33	0.41	0	44,47,47	0.85	2 (4%)
13	NAG	C	804	1	14,14,15	0.33	0	17,19,21	0.65	0
13	NAG	E	803	1	14,14,15	0.37	0	17,19,21	1.25	3 (17%)
13	NAG	B	801	1	14,14,15	0.37	0	17,19,21	0.53	0
13	NAG	E	802	1	14,14,15	0.42	0	17,19,21	0.75	0
13	NAG	C	803	1	14,14,15	0.23	0	17,19,21	0.96	2 (11%)
13	NAG	A	802	1	14,14,15	0.28	0	17,19,21	0.88	1 (5%)
13	NAG	A	804	1	14,14,15	0.38	0	17,19,21	0.57	0
13	NAG	D	801	1	14,14,15	0.40	0	17,19,21	1.01	2 (11%)

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
13	NAG	A	803	1	-	1/6/23/26	0/1/1/1
13	NAG	C	802	1	-	0/6/23/26	0/1/1/1
13	NAG	C	805	1	-	2/6/23/26	0/1/1/1
14	83G	E	801	-	-	6/22/35/35	0/4/4/4
14	83G	A	801	-	-	13/22/35/35	0/4/4/4
14	83G	C	801	-	-	8/22/35/35	0/4/4/4
13	NAG	C	804	1	-	2/6/23/26	0/1/1/1
13	NAG	E	803	1	-	3/6/23/26	0/1/1/1
13	NAG	B	801	1	-	0/6/23/26	0/1/1/1
13	NAG	E	802	1	-	1/6/23/26	0/1/1/1
13	NAG	C	803	1	-	5/6/23/26	0/1/1/1
13	NAG	A	802	1	-	3/6/23/26	0/1/1/1
13	NAG	A	804	1	-	0/6/23/26	0/1/1/1
13	NAG	D	801	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	D	801	NAG	C1-C2-N2	2.93	115.05	110.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	E	801	83G	C11-C12-C14	2.75	124.22	118.82
13	C	805	NAG	O4-C4-C5	2.59	115.71	109.32
14	C	801	83G	O02-C03-C08	2.53	119.42	115.84
13	E	803	NAG	C1-O5-C5	2.47	115.50	112.19
13	A	802	NAG	C2-N2-C7	2.44	126.17	122.90
14	E	801	83G	O02-C03-C08	2.42	119.27	115.84
13	D	801	NAG	C2-N2-C7	2.41	126.13	122.90
14	A	801	83G	C12-C14-N16	2.22	121.08	118.81
13	E	803	NAG	C2-N2-C7	2.21	125.87	122.90
13	C	803	NAG	C2-N2-C7	2.20	125.85	122.90
14	C	801	83G	C11-C12-C14	2.09	122.92	118.82
14	E	801	83G	C21-N16-C14	2.08	125.50	120.32
13	E	803	NAG	O4-C4-C5	2.05	114.38	109.32
13	C	803	NAG	C4-C3-C2	-2.04	108.03	111.02

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
13	C	804	NAG	C8-C7-N2-C2
13	C	804	NAG	O7-C7-N2-C2
13	D	801	NAG	C8-C7-N2-C2
13	D	801	NAG	O7-C7-N2-C2
13	E	803	NAG	C1-C2-N2-C7
13	E	803	NAG	C8-C7-N2-C2
13	E	803	NAG	O7-C7-N2-C2
13	A	802	NAG	C1-C2-N2-C7
13	A	802	NAG	C8-C7-N2-C2
13	A	802	NAG	O7-C7-N2-C2
14	C	801	83G	O15-C14-N16-C17
14	C	801	83G	C08-C11-C12-O13
14	A	801	83G	O15-C14-N16-C17
14	A	801	83G	C10-C11-C12-O13
14	A	801	83G	C08-C11-C12-C14
14	A	801	83G	C08-C11-C12-O13
13	C	803	NAG	O5-C5-C6-O6
13	C	803	NAG	C8-C7-N2-C2
13	C	803	NAG	O7-C7-N2-C2
14	A	801	83G	C25-C23-N19-C20
13	C	805	NAG	O5-C5-C6-O6
14	A	801	83G	O24-C23-N19-C20
14	A	801	83G	C04-C03-O02-C01

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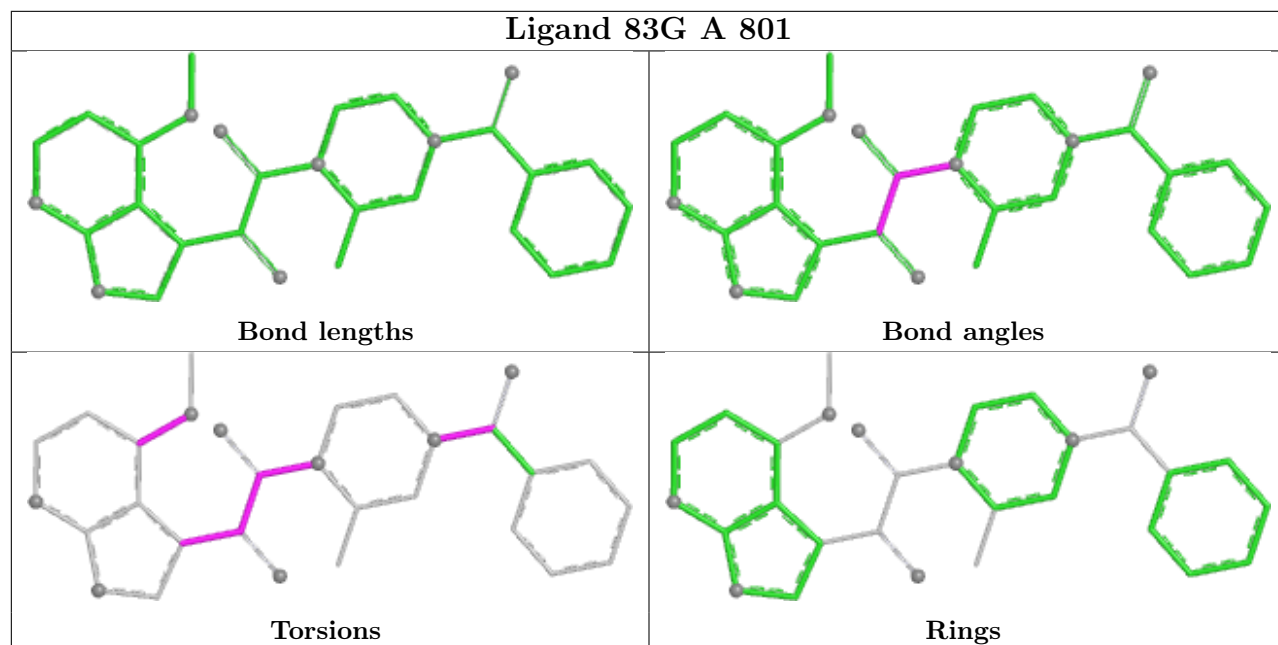
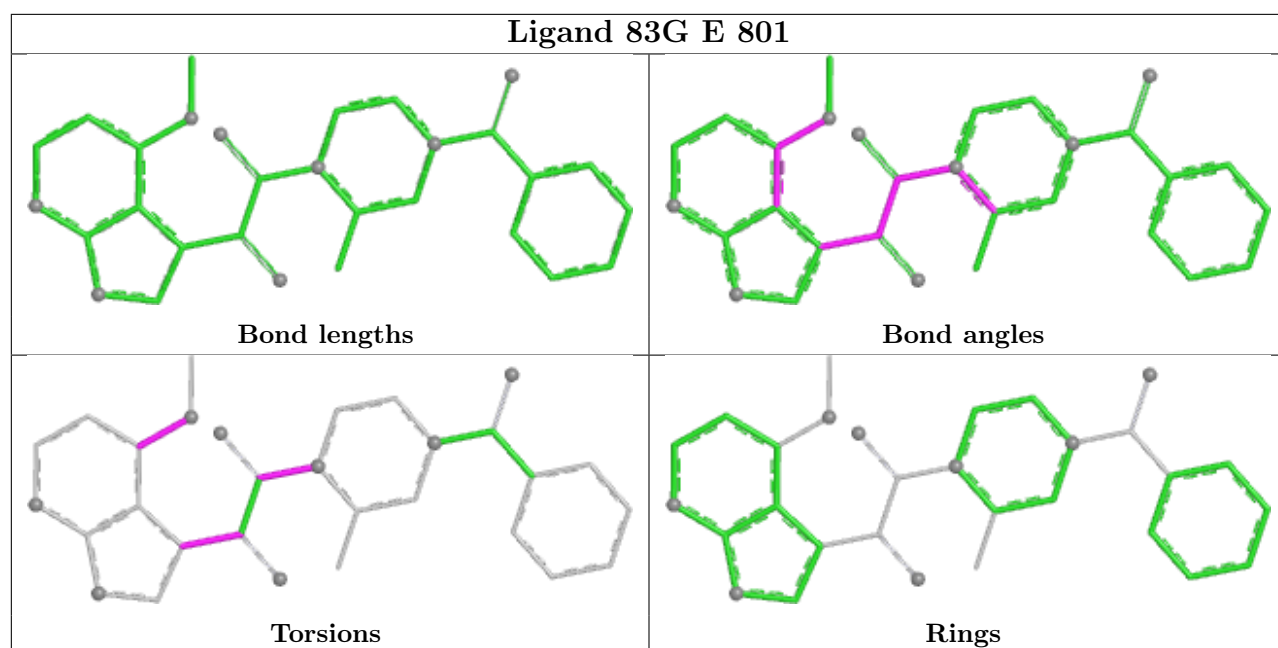
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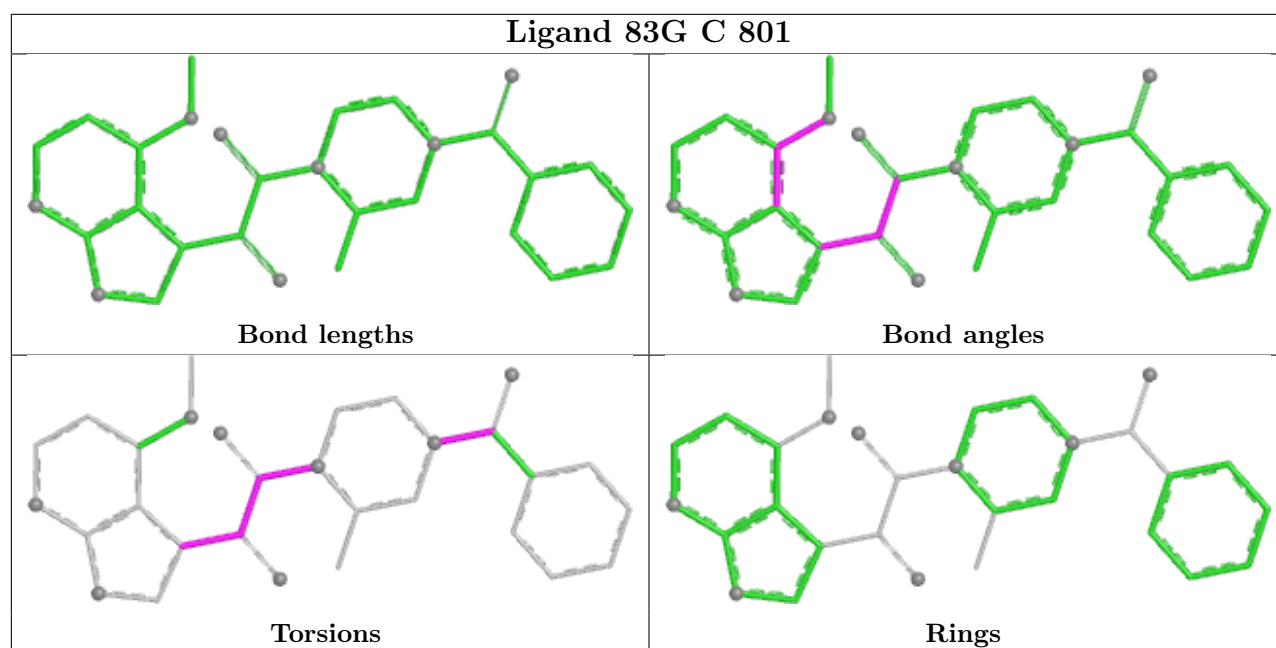
Mol	Chain	Res	Type	Atoms
13	C	803	NAG	C4-C5-C6-O6
14	A	801	83G	C08-C03-O02-C01
14	C	801	83G	C12-C14-N16-C17
14	E	801	83G	O15-C14-N16-C17
13	C	803	NAG	C3-C2-N2-C7
13	D	801	NAG	C3-C2-N2-C7
14	A	801	83G	C12-C14-N16-C17
14	E	801	83G	C04-C03-O02-C01
14	E	801	83G	C08-C03-O02-C01
13	D	801	NAG	C1-C2-N2-C7
14	A	801	83G	C11-C12-C14-O15
14	A	801	83G	C10-C11-C12-C14
13	C	805	NAG	C4-C5-C6-O6
14	E	801	83G	C08-C11-C12-O13
14	C	801	83G	C10-C11-C12-O13
14	E	801	83G	C10-C11-C12-O13
14	C	801	83G	C12-C14-N16-C21
14	E	801	83G	C12-C14-N16-C21
14	A	801	83G	C12-C14-N16-C21
14	C	801	83G	O24-C23-N19-C20
13	E	802	NAG	C1-C2-N2-C7
13	A	803	NAG	C1-C2-N2-C7
14	C	801	83G	C11-C12-C14-O15
14	C	801	83G	O13-C12-C14-O15
14	A	801	83G	O13-C12-C14-O15

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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.

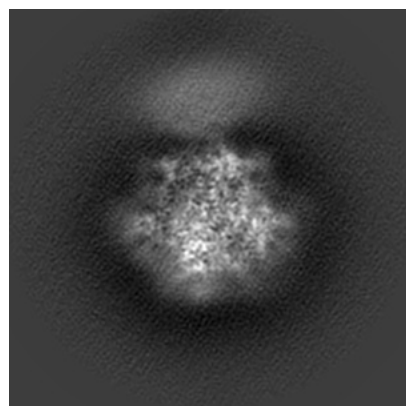
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-61554. These allow visual inspection of the internal detail of the map and identification of artifacts.

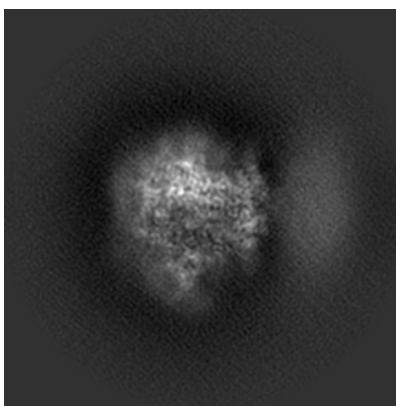
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

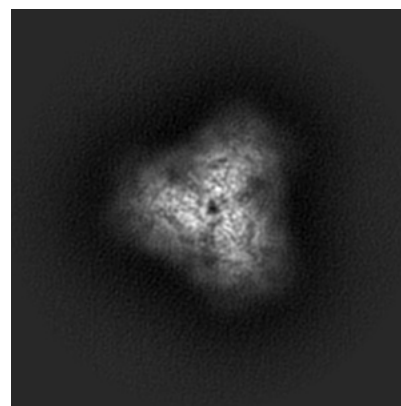
6.1.1 Primary map



X

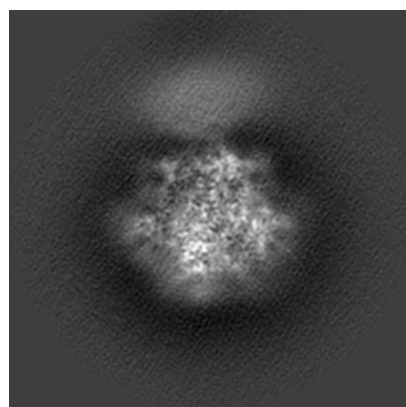


Y

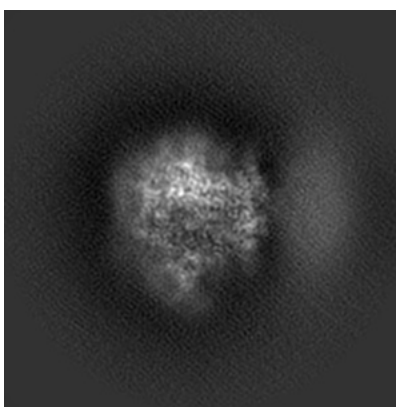


Z

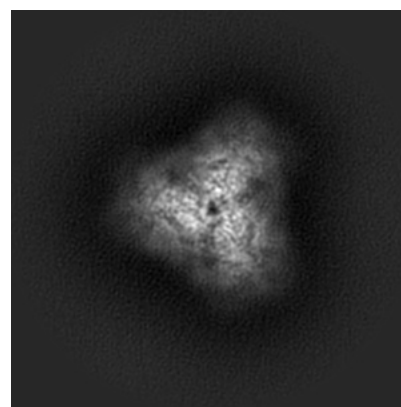
6.1.2 Raw map



X



Y

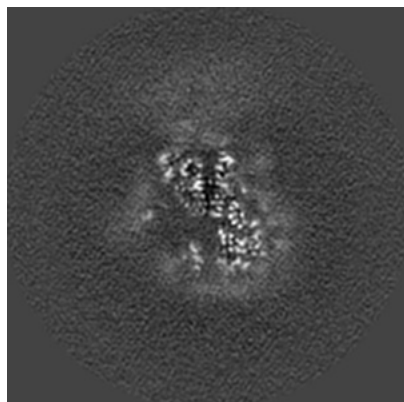


Z

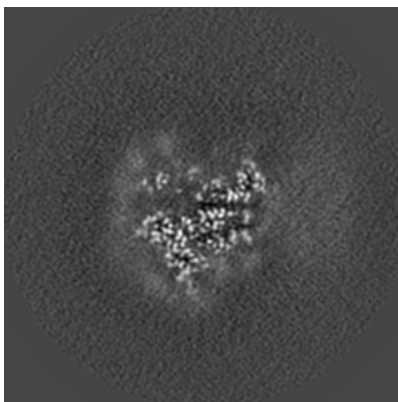
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

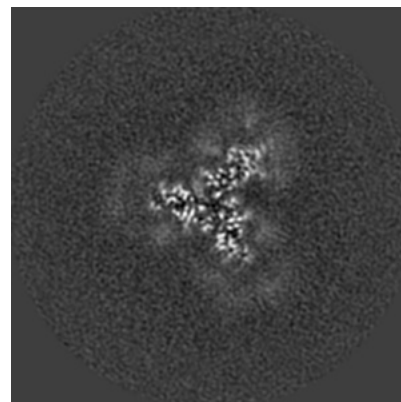
6.2.1 Primary map



X Index: 90

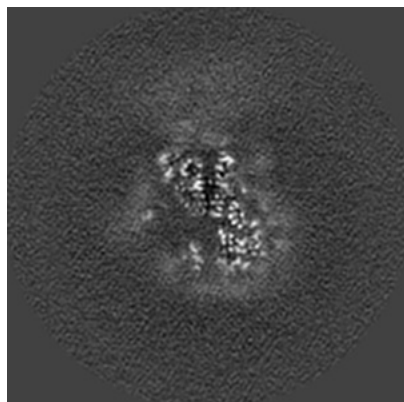


Y Index: 90

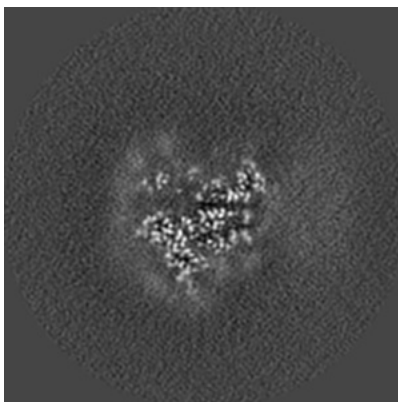


Z Index: 90

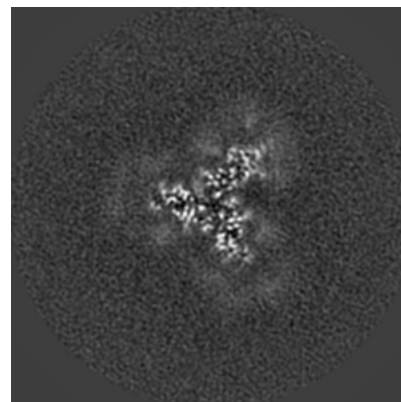
6.2.2 Raw map



X Index: 90



Y Index: 90

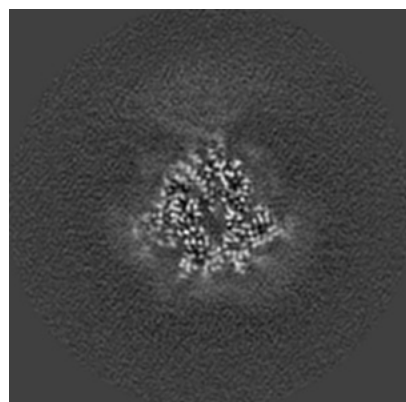


Z Index: 90

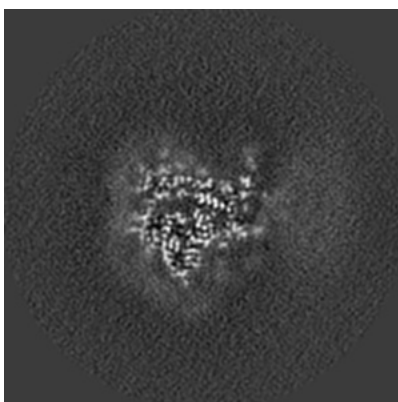
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

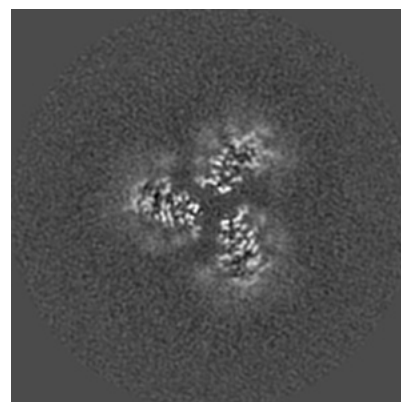
6.3.1 Primary map



X Index: 98

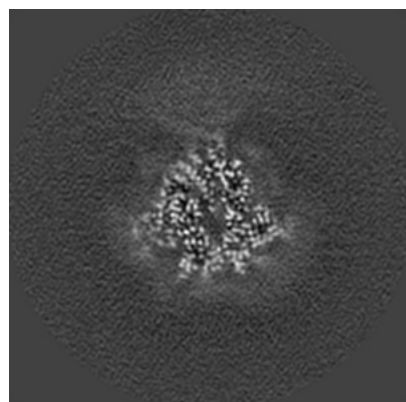


Y Index: 87

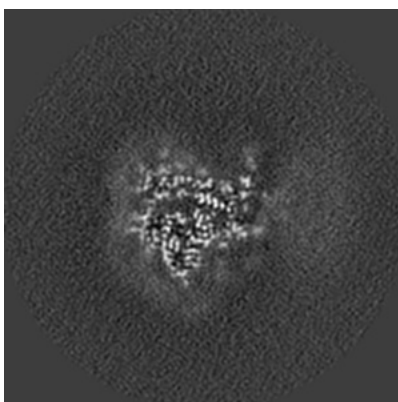


Z Index: 83

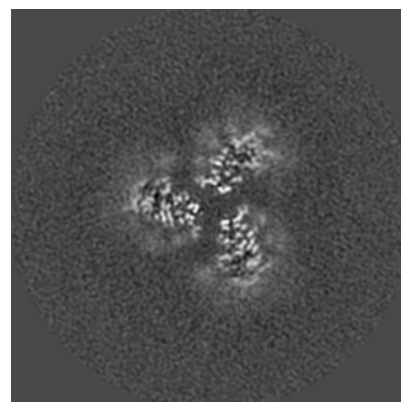
6.3.2 Raw map



X Index: 98



Y Index: 87

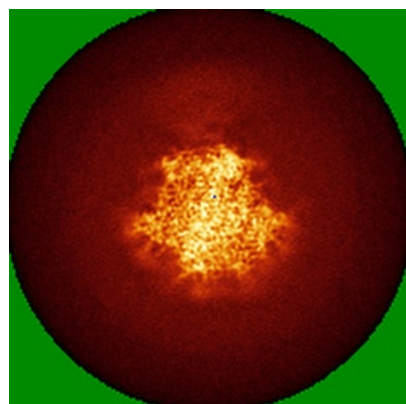


Z Index: 83

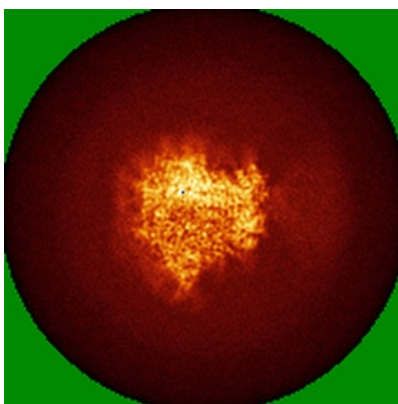
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

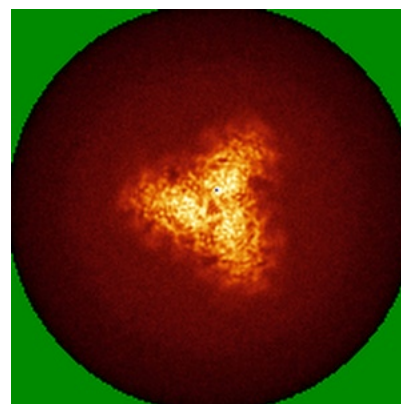
6.4.1 Primary map



X

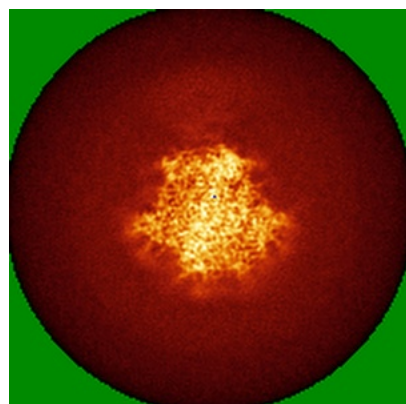


Y

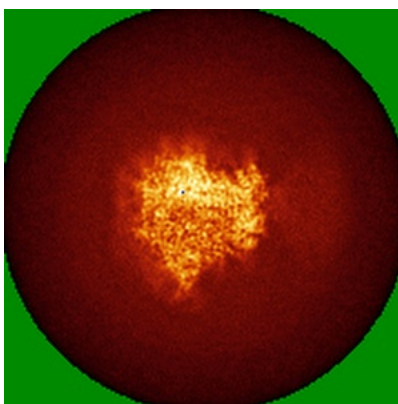


Z

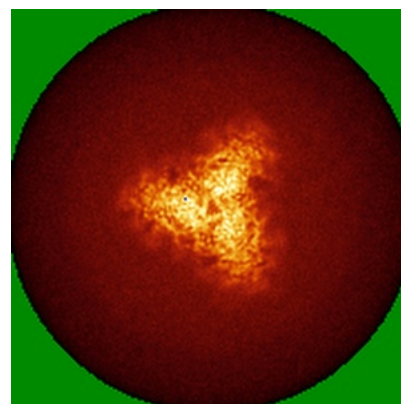
6.4.2 Raw map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.02. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

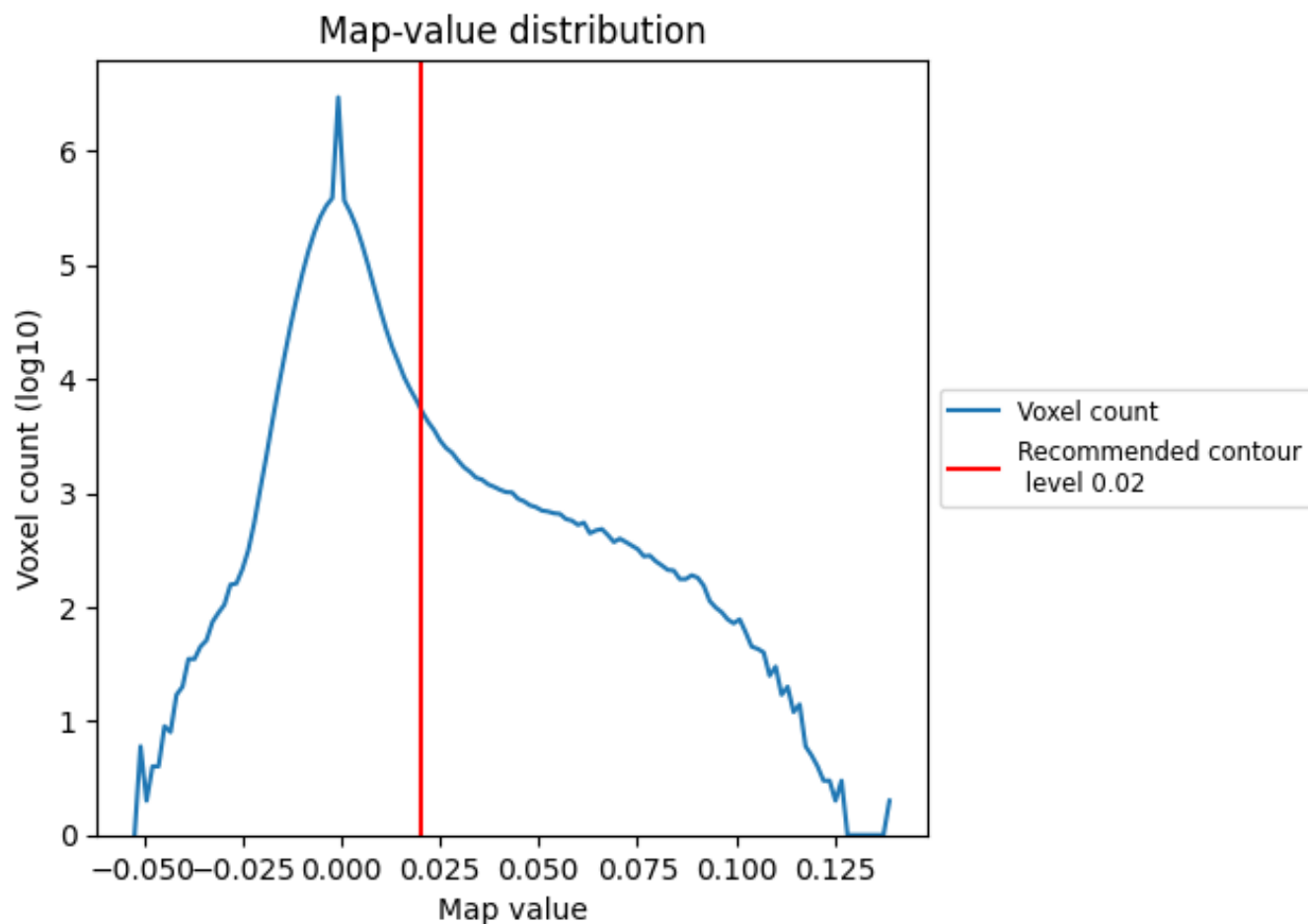
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

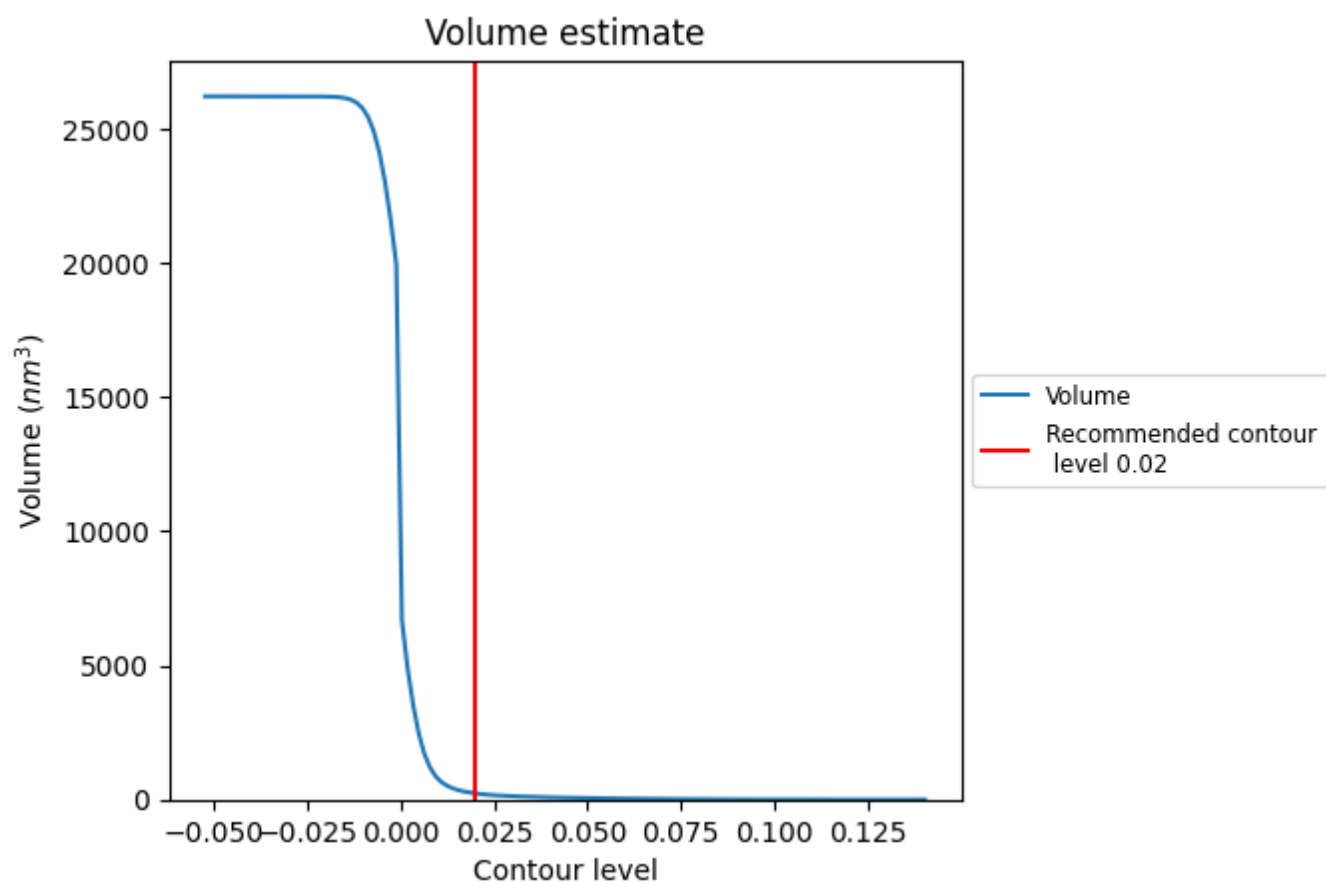
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

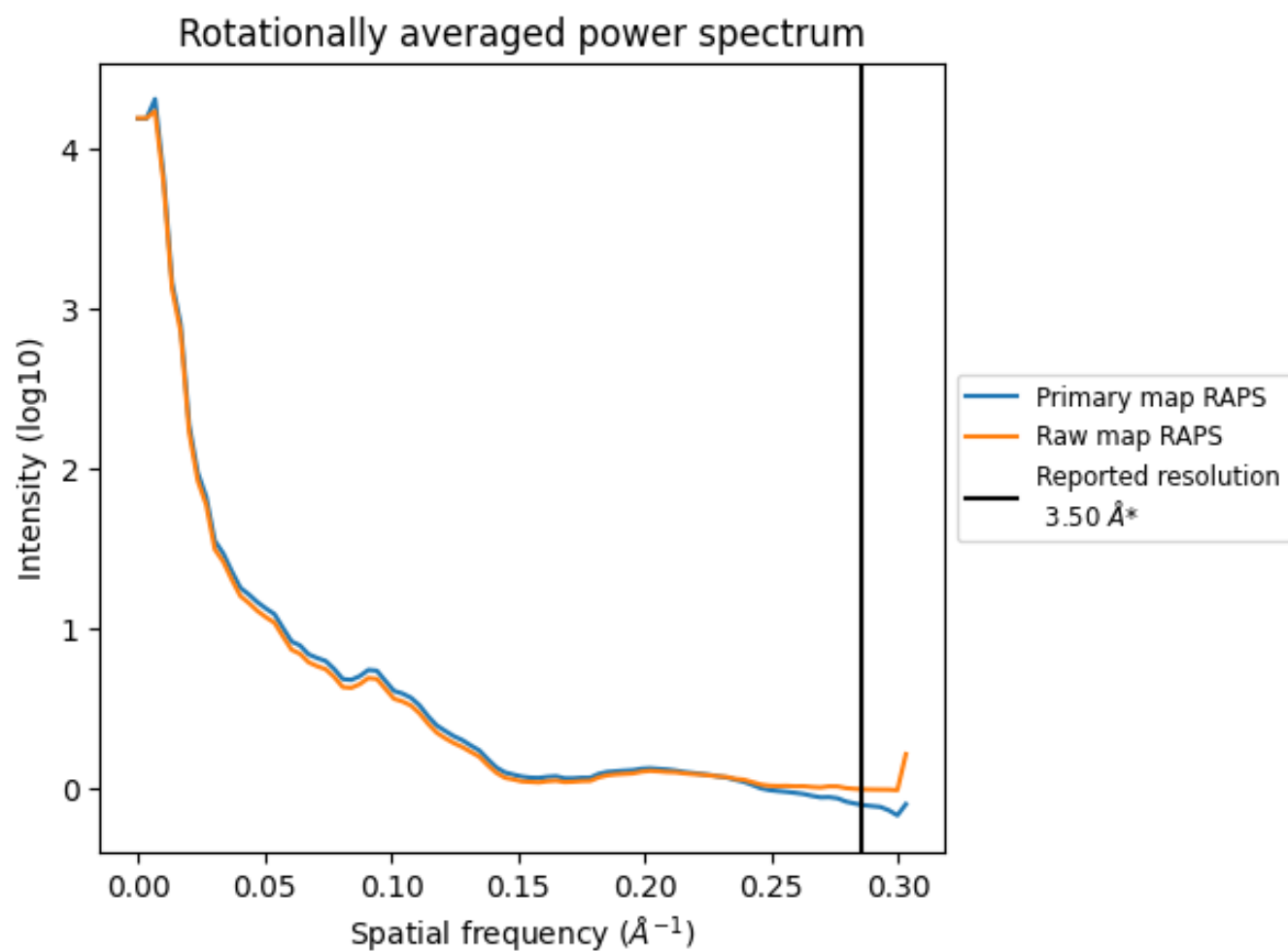
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 229 nm^3 ; this corresponds to an approximate mass of 207 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

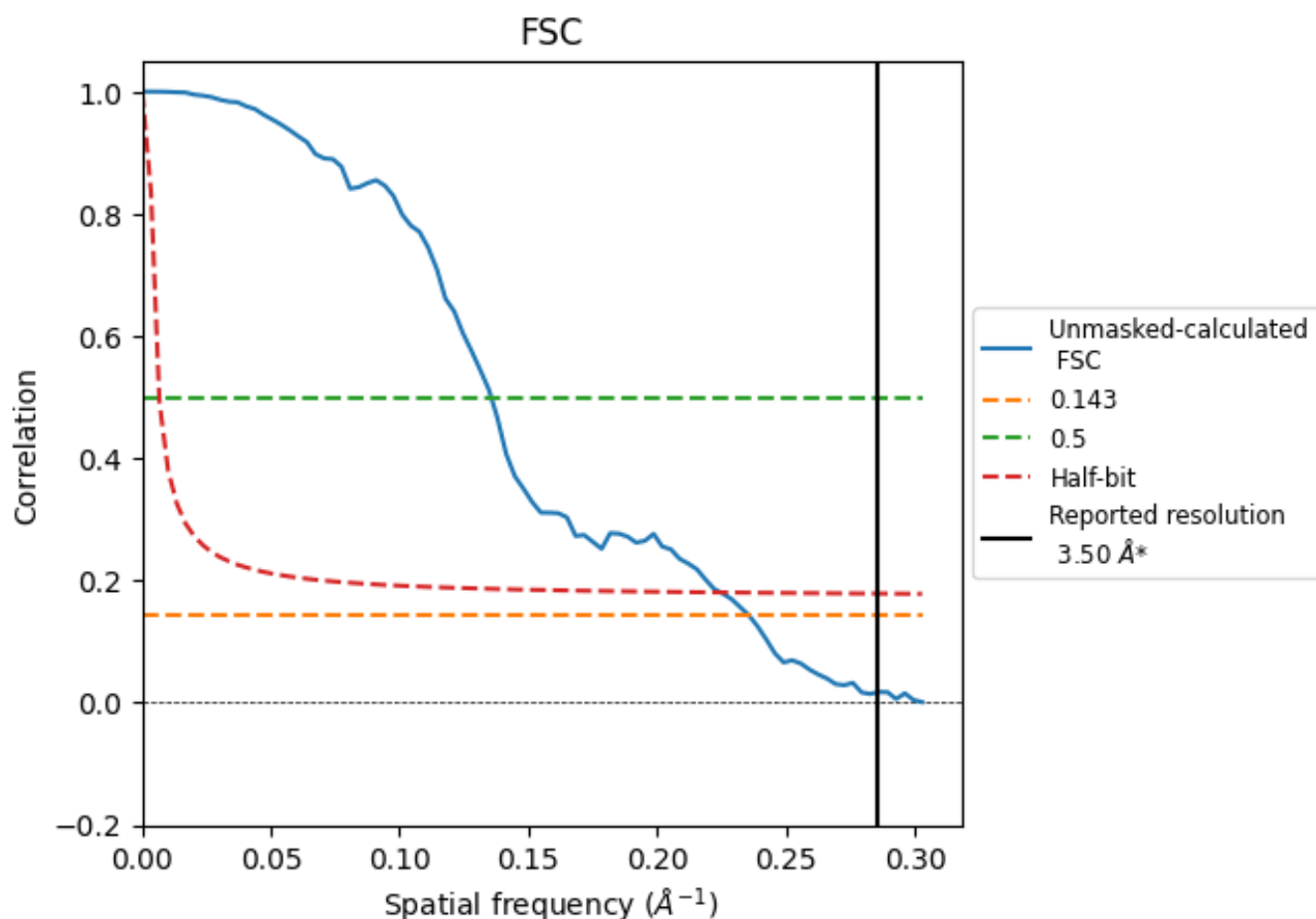


*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.286 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

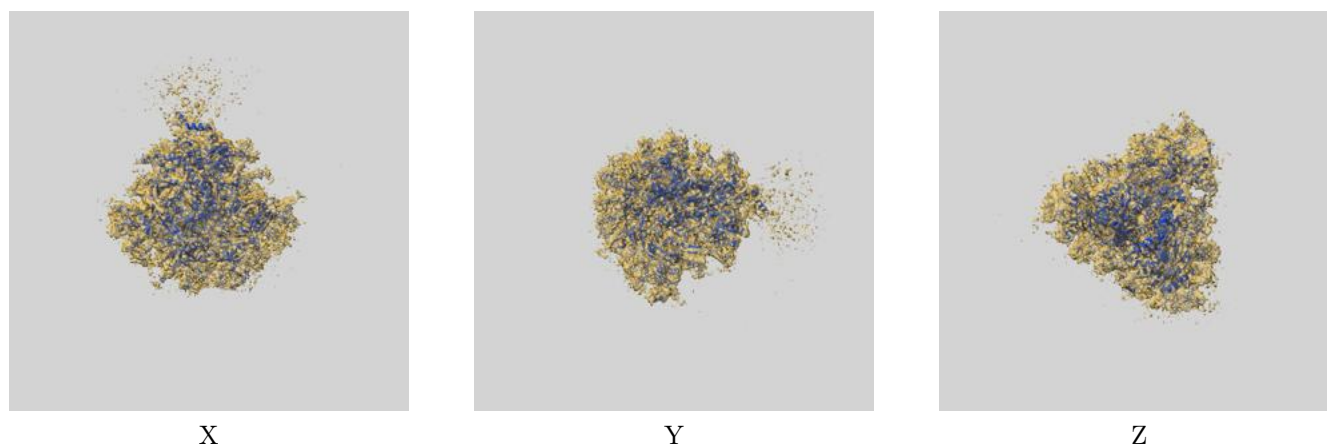
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.24	7.37	4.45

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.24 differs from the reported value 3.5 by more than 10 %

9 Map-model fit [i](#)

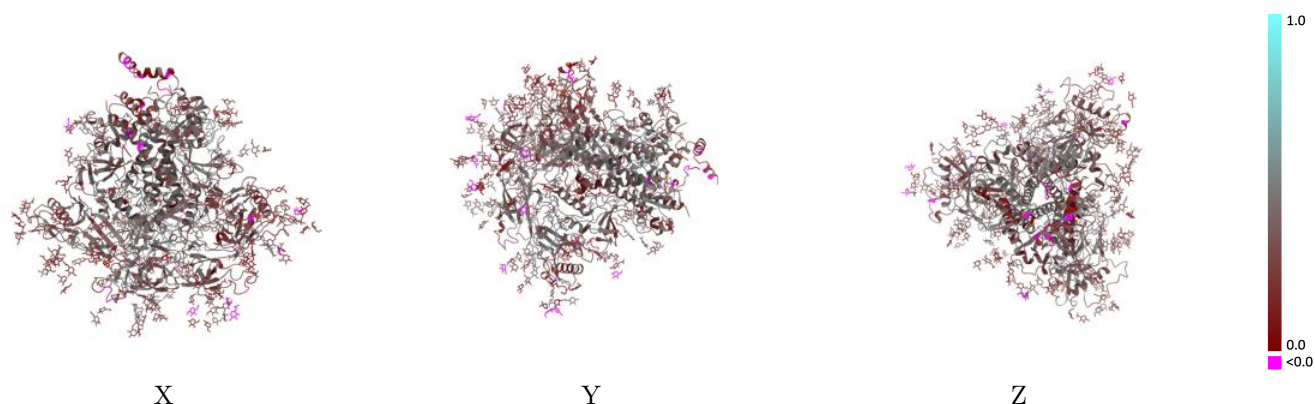
This section contains information regarding the fit between EMDB map EMD-61554 and PDB model 9JKG. Per-residue inclusion information can be found in section 3 on page 19.

9.1 Map-model overlay [i](#)



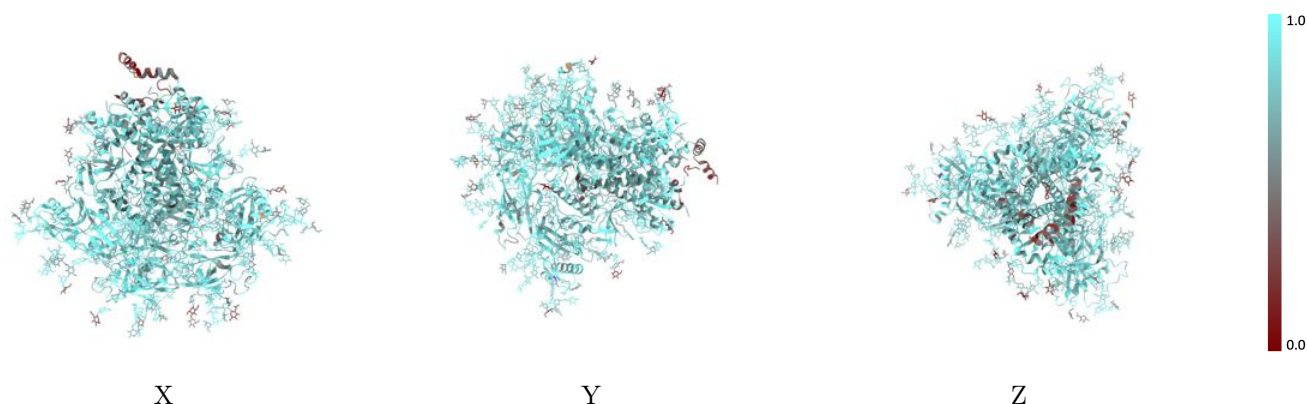
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



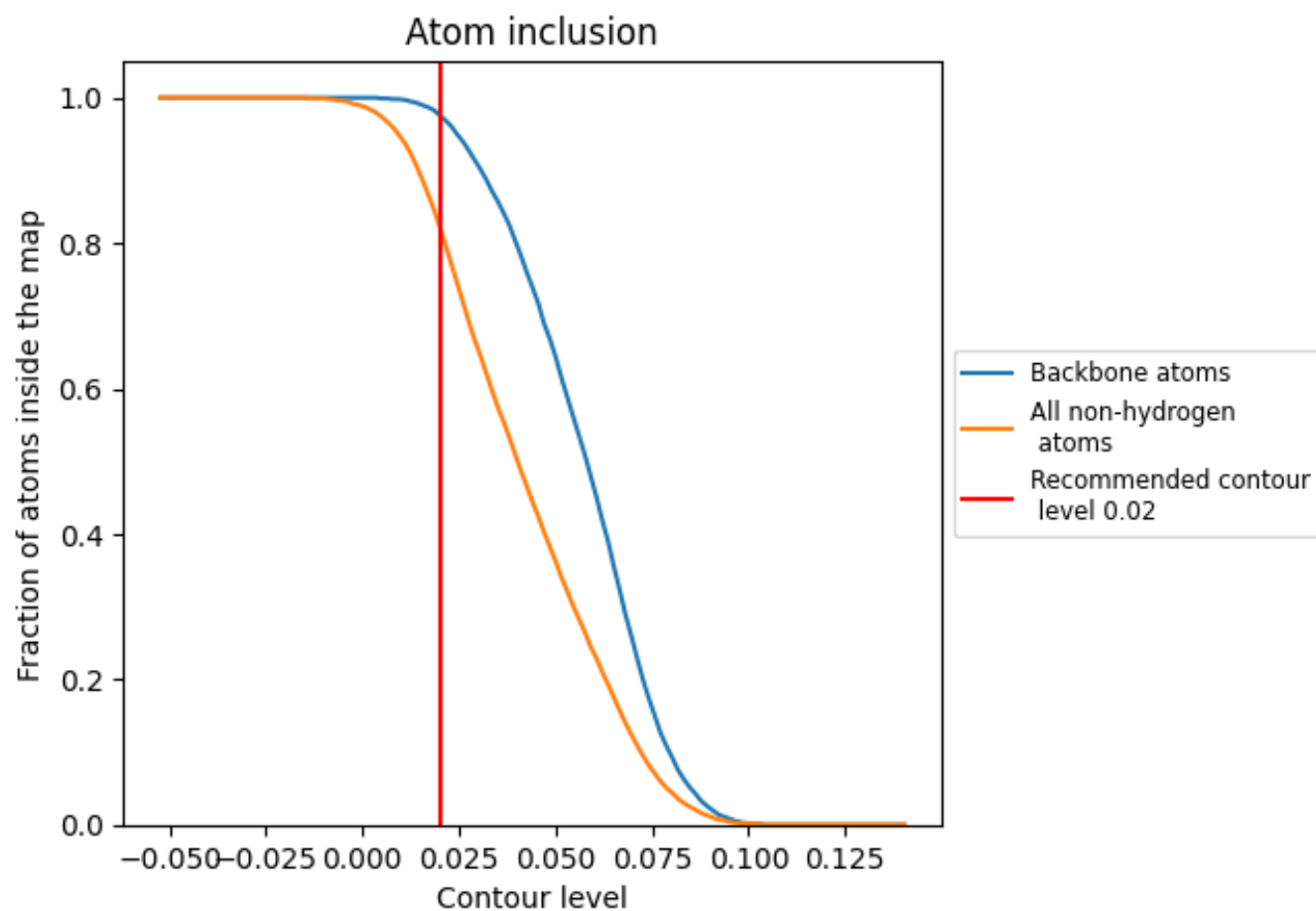
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.02).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 82% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8230	 0.3380
0	 0.6920	 0.2400
1	 0.6150	 0.1720
2	 0.9310	 0.3280
3	 0.7400	 0.1900
4	 0.7180	 0.1950
5	 0.7200	 0.2870
6	 0.8690	 0.3490
7	 0.7230	 0.1370
8	 0.8460	 0.3100
9	 0.8000	 0.2470
A	 0.8580	 0.3670
AA	 0.8720	 0.2720
B	 0.7340	 0.3230
BA	 0.7440	 0.2390
C	 0.8540	 0.3650
CA	 0.7140	 0.2450
D	 0.7990	 0.3460
DA	 0.3680	 0.2500
E	 0.8560	 0.3440
EA	 0.5360	 0.2090
F	 0.8040	 0.3660
FA	 0.5530	 0.2210
G	 0.8970	 0.3320
H	 0.4640	 0.1490
I	 0.7950	 0.3490
J	 0.7860	 0.3620
K	 0.8400	 0.3120
L	 0.8200	 0.3100
M	 0.6070	 0.2860
N	 0.8720	 0.2830
O	 0.4290	 0.1040
P	 0.7140	 0.2800
Q	 0.8890	 0.3830
R	 0.7440	 0.2460



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Chain	Atom inclusion	Q-score
S	 0.7860	 0.2740
T	 0.6920	 0.1910
U	 0.7600	 0.1320
V	 0.9180	 0.2440
W	 0.7800	 0.1230
X	 0.7400	 0.2070
Y	 0.7950	 0.2680
Z	 0.7180	 0.2600
a	 0.6730	 0.2440
b	 0.3930	 0.1140
c	 0.6670	 0.2680
d	 0.7950	 0.2790
e	 0.8530	 0.3050
f	 0.7380	 0.1120
g	 0.4870	 0.1150
h	 0.5000	 0.0910
i	 0.8000	 0.2200
j	 0.5900	 0.1890
k	 0.7180	 0.3020
l	 0.7770	 0.3190
m	 0.8200	 0.2490
n	 0.6670	 0.3100
o	 0.8210	 0.1960
p	 0.8600	 0.3140
q	 0.7830	 0.1240
r	 0.7600	 0.1730
s	 0.6920	 0.1920
t	 0.7950	 0.2410
u	 0.8210	 0.3010
v	 0.7500	 0.3340
w	 0.8360	 0.3500
x	 0.8460	 0.2740
y	 0.8460	 0.2440
z	 0.6920	 0.2420