



Full wwPDB EM Validation Report ⓘ

Mar 23, 2026 – 03:33 PM UTC

PDB ID : 6CUE / pdb_00006cue
EMDB ID : EMD-7621
Title : Cryo-EM structure at 4.0 Å resolution of vaccine-elicited antibody vFP7.04 in complex with HIV-1 Env BG505 DS-SOSIP, and antibodies VRC03 and PGT122
Authors : Acharya, P.; Carragher, B.; Potter, C.S.; Kwong, P.D.
Deposited on : 2018-03-26
Resolution : 4.00 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:

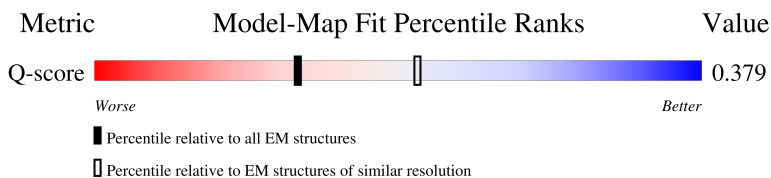
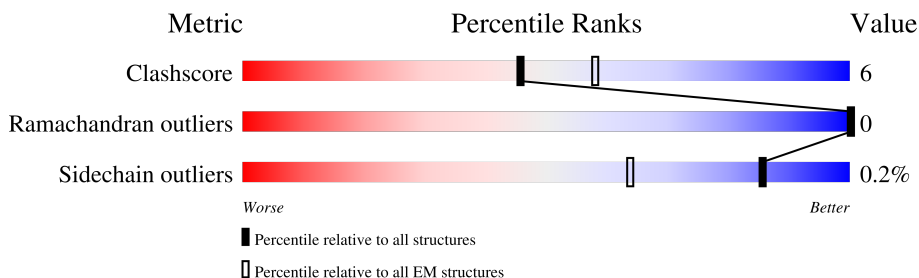
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 4.00 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.








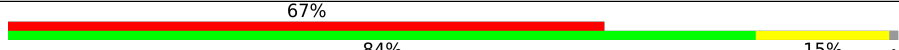


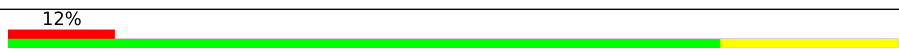

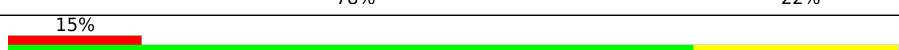
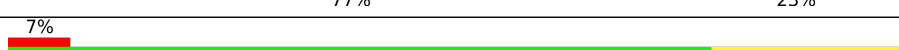

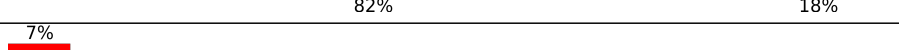
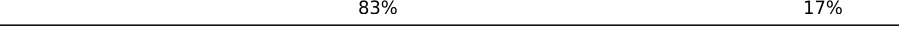
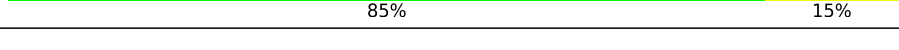







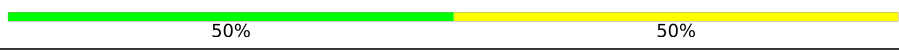
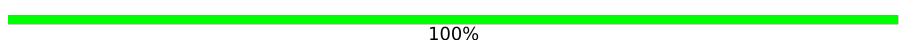
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	7587 (3.50 - 4.50)

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	2	473	<p>82% 14% .</p>
1	C	473	<p>82% 13% .</p>
1	c	473	<p>80% 16% .</p>
2	1	153	<p>74% 12% 14%</p>


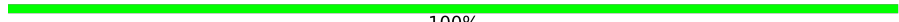






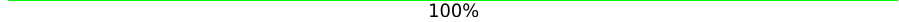



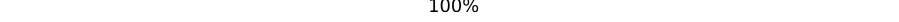
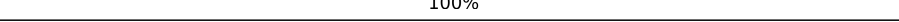
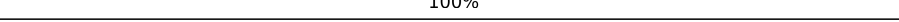

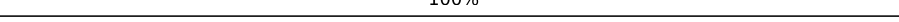
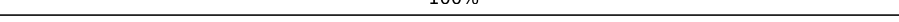







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Mol	Chain	Length	Quality of chain
2	D	153	
2	d	153	
3	3	118	
3	H	118	
3	h	118	
4	4	112	
4	L	112	
4	l	112	
5	5	132	
5	M	132	
5	m	132	
6	6	107	
6	N	107	
6	n	107	
7	7	129	
7	Q	129	
7	q	129	
8	8	102	
8	R	102	
8	r	102	
9	9	2	
9	A	2	
9	F	2	
9	G	2	
9	J	2	




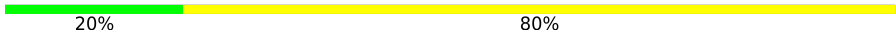
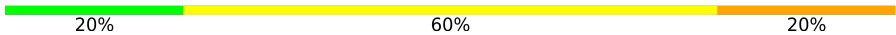
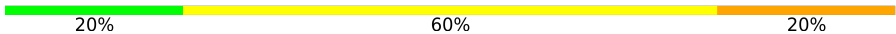
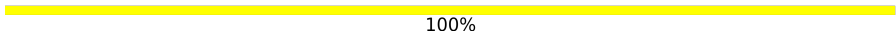
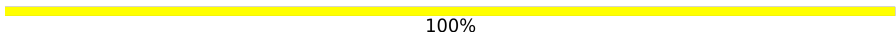
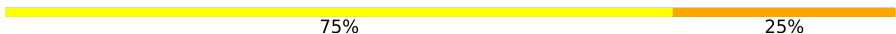
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Mol	Chain	Length	Quality of chain
9	K	2	 100%
9	P	2	 100%
9	T	2	 50% 50%
9	V	2	 50% 50%
9	Y	2	 50% 50%
9	Z	2	 100%
9	b	2	 100%
9	e	2	 50% 50%
9	g	2	 100%
9	j	2	 50% 50%
9	o	2	 50% 50%
9	t	2	 50% 50%
9	u	2	 100%
9	w	2	 100%
9	x	2	 100%
9	z	2	 50% 50%
10	B	3	 100%
10	E	3	 100%
10	O	3	 67% 33%
10	W	3	 100%
10	X	3	 100%
10	f	3	 67% 33%
10	p	3	 100%
10	s	3	 100%
10	y	3	 67% 33%

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Mol	Chain	Length	Quality of chain
11	I	5	 40% 60%
11	a	5	 60% 40%
11	v	5	 60% 40%
12	0	5	 20% 80%
12	S	5	 20% 60% 20%
12	i	5	 20% 60% 20%
13	AA	4	 100%
13	U	4	 100%
13	k	4	 75% 25%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 31692 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 gp120.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	c	453	3564	2233	630	671	30	0	0
1	2	453	3564	2233	630	671	30	0	0
1	C	453	3564	2233	630	671	30	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
c	201	CYS	ILE	conflict	UNP Q2N0S6
c	332	ASN	THR	conflict	UNP Q2N0S6
c	433	CYS	ALA	conflict	UNP Q2N0S6
c	501	CYS	ALA	conflict	UNP Q2N0S6
2	201	CYS	ILE	conflict	UNP Q2N0S6
2	332	ASN	THR	conflict	UNP Q2N0S6
2	433	CYS	ALA	conflict	UNP Q2N0S6
2	501	CYS	ALA	conflict	UNP Q2N0S6
C	201	CYS	ILE	conflict	UNP Q2N0S6
C	332	ASN	THR	conflict	UNP Q2N0S6
C	433	CYS	ALA	conflict	UNP Q2N0S6
C	501	CYS	ALA	conflict	UNP Q2N0S6

- Molecule 2 is a protein called Envelope glycoprotein gp41.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	1	132	1034	654	178	196	6	0	0
2	D	132	1034	654	178	196	6	0	0
2	d	132	1034	654	178	196	6	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	605	CYS	THR	conflict	UNP Q2N0S7
D	605	CYS	THR	conflict	UNP Q2N0S7
d	605	CYS	THR	conflict	UNP Q2N0S7

- Molecule 3 is a protein called vFP7.04 Heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	117	Total	C	N	O	S	0	0
			919	589	149	176	5		
3	H	117	Total	C	N	O	S	0	0
			919	589	149	176	5		
3	h	117	Total	C	N	O	S	0	0
			919	589	149	176	5		

- Molecule 4 is a protein called vFP7.04 light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	4	111	Total	C	N	O	S	0	0
			853	540	145	165	3		
4	L	111	Total	C	N	O	S	0	0
			853	540	145	165	3		
4	l	111	Total	C	N	O	S	0	0
			853	540	145	165	3		

- Molecule 5 is a protein called PGT122 heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	5	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		
5	M	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		
5	m	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		

- Molecule 6 is a protein called PGT122 Light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	6	107	Total	C	N	O	S	0	0
			817	512	141	162	2		
6	N	107	Total	C	N	O	S	0	0
			817	512	141	162	2		

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Mol	Chain	Residues	Atoms					AltConf	Trace
6	n	107	Total	C	N	O	S	0	0
			817	512	141	162	2		

- Molecule 7 is a protein called VRC03 Heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	7	129	Total	C	N	O	S	0	0
			1029	660	176	187	6		
7	Q	129	Total	C	N	O	S	0	0
			1029	660	176	187	6		
7	q	129	Total	C	N	O	S	0	0
			1029	660	176	187	6		

- Molecule 8 is a protein called VRC03 light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	8	102	Total	C	N	O	S	0	0
			802	510	137	152	3		
8	R	102	Total	C	N	O	S	0	0
			802	510	137	152	3		
8	r	102	Total	C	N	O	S	0	0
			802	510	137	152	3		

- Molecule 9 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
9	A	2	Total	C	N	O	0	0
			28	16	2	10		
9	F	2	Total	C	N	O	0	0
			28	16	2	10		
9	G	2	Total	C	N	O	0	0
			28	16	2	10		
9	J	2	Total	C	N	O	0	0
			28	16	2	10		
9	K	2	Total	C	N	O	0	0
			28	16	2	10		

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Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
9	P	2	28	16	2	10	0	0
9	T	2	28	16	2	10	0	0
9	V	2	28	16	2	10	0	0
9	Y	2	28	16	2	10	0	0
9	Z	2	28	16	2	10	0	0
9	b	2	28	16	2	10	0	0
9	e	2	28	16	2	10	0	0
9	g	2	28	16	2	10	0	0
9	j	2	28	16	2	10	0	0
9	o	2	28	16	2	10	0	0
9	t	2	28	16	2	10	0	0
9	u	2	28	16	2	10	0	0
9	w	2	28	16	2	10	0	0
9	x	2	28	16	2	10	0	0
9	z	2	28	16	2	10	0	0
9	9	2	28	16	2	10	0	0

- Molecule 10 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.



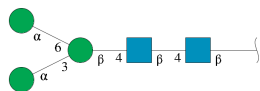
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
10	B	3	39	22	2	15	0	0

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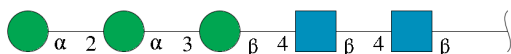
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
10	E	3	Total 39	C 22	N 2	O 15	0	0
10	O	3	Total 39	C 22	N 2	O 15	0	0
10	W	3	Total 39	C 22	N 2	O 15	0	0
10	X	3	Total 39	C 22	N 2	O 15	0	0
10	f	3	Total 39	C 22	N 2	O 15	0	0
10	p	3	Total 39	C 22	N 2	O 15	0	0
10	s	3	Total 39	C 22	N 2	O 15	0	0
10	y	3	Total 39	C 22	N 2	O 15	0	0

- Molecule 11 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
			Total	C	N	O		
11	I	5	Total 61	C 34	N 2	O 25	0	0
11	a	5	Total 61	C 34	N 2	O 25	0	0
11	v	5	Total 61	C 34	N 2	O 25	0	0

- Molecule 12 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-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.



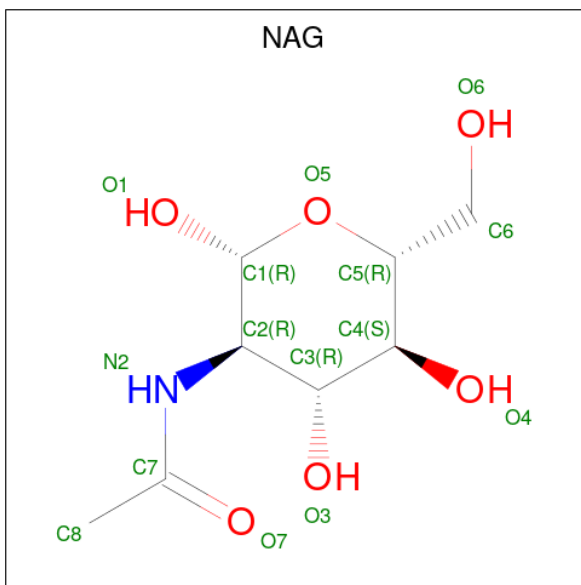
Mol	Chain	Residues	Atoms				AltConf	Trace
12	S	5	Total	C	N	O	0	0
			61	34	2	25		
12	i	5	Total	C	N	O	0	0
			61	34	2	25		
12	0	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 13 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
13	U	4	Total	C	N	O	0	0
			50	28	2	20		
13	k	4	Total	C	N	O	0	0
			50	28	2	20		
13	AA	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 14 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

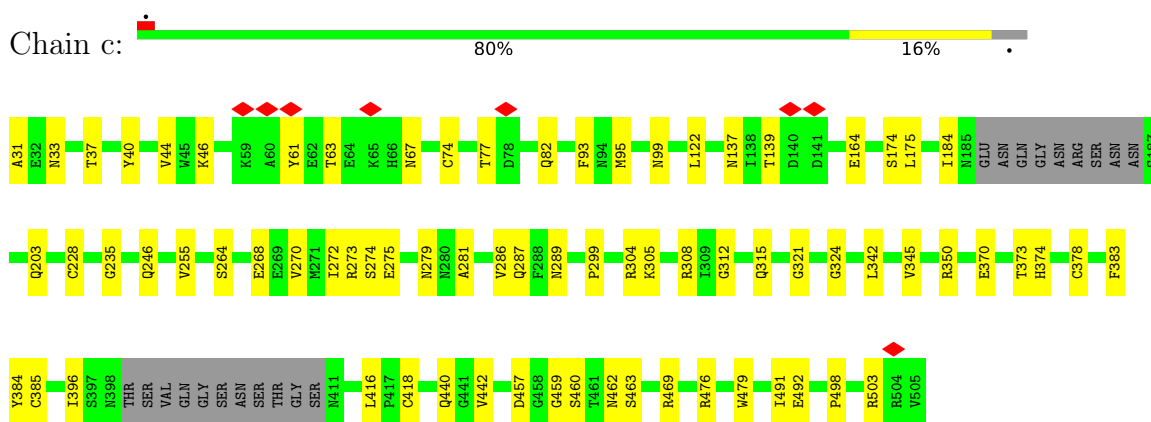


Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
14	c	1	14	8	1	5	0
14	2	1	14	8	1	5	0
14	C	1	14	8	1	5	0

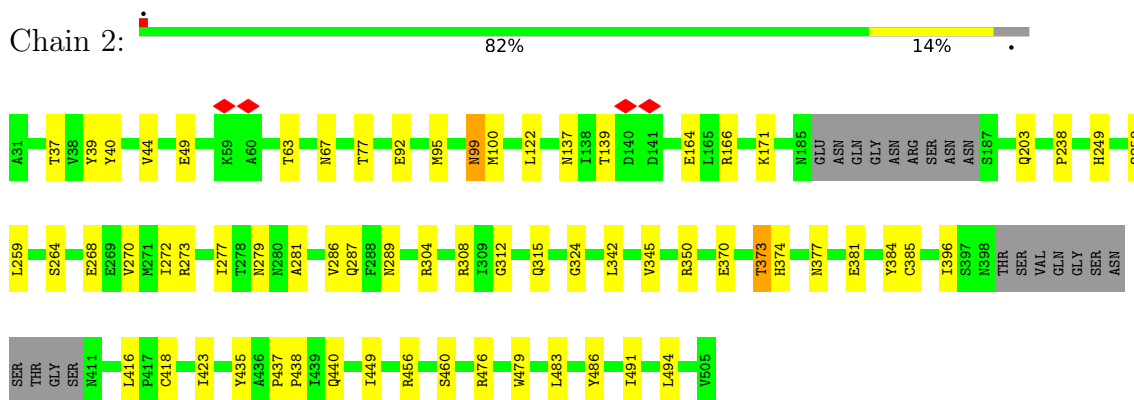
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

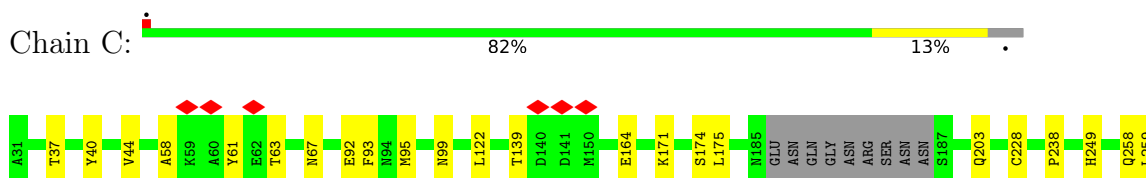
- Molecule 1: Envelope glycoprotein gp120

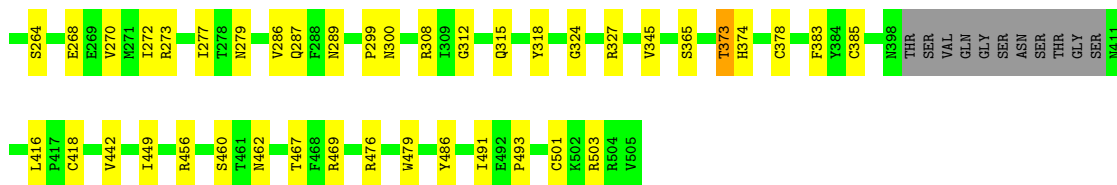


- Molecule 1: Envelope glycoprotein gp120

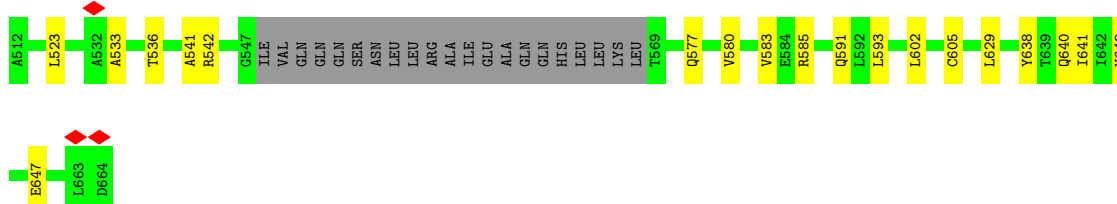


- Molecule 1: Envelope glycoprotein gp120

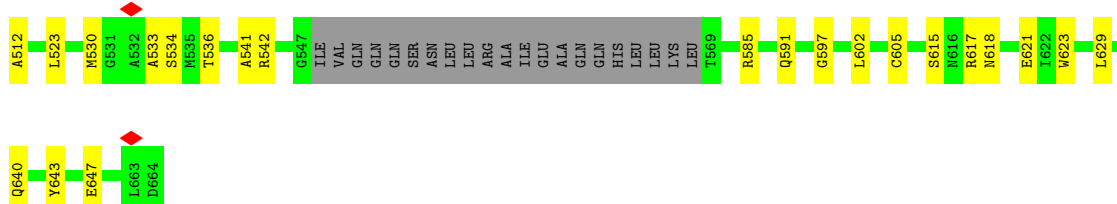




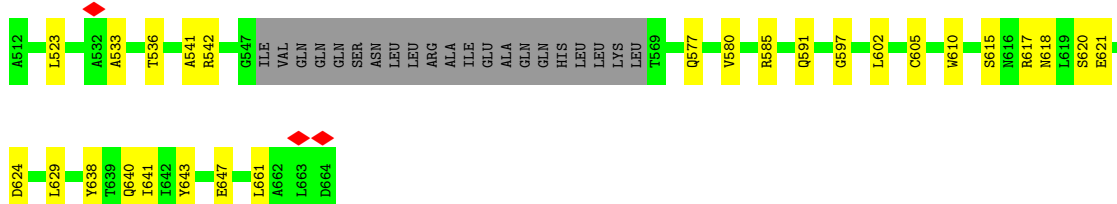
• Molecule 2: Envelope glycoprotein gp41



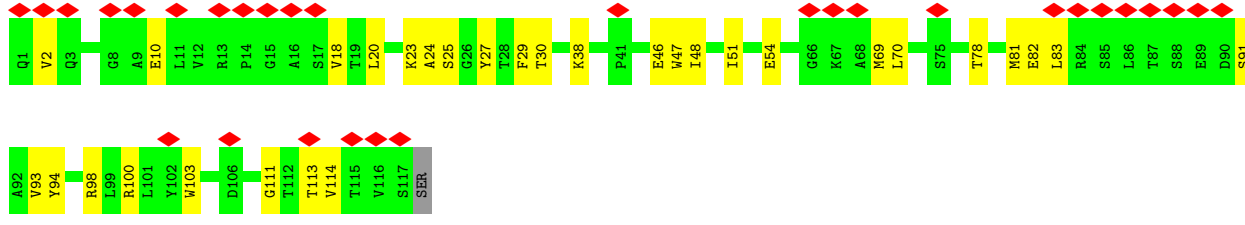
• Molecule 2: Envelope glycoprotein gp41



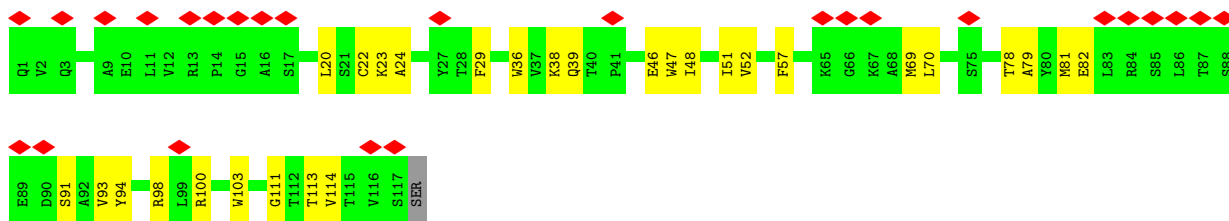
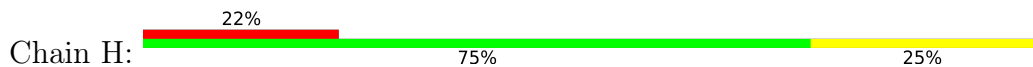
• Molecule 2: Envelope glycoprotein gp41



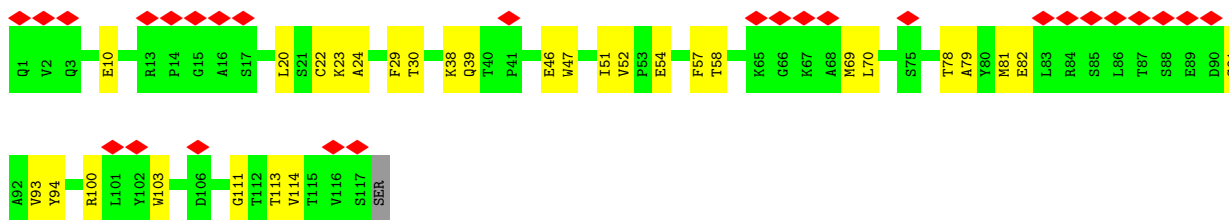
• Molecule 3: vFP7.04 Heavy chain



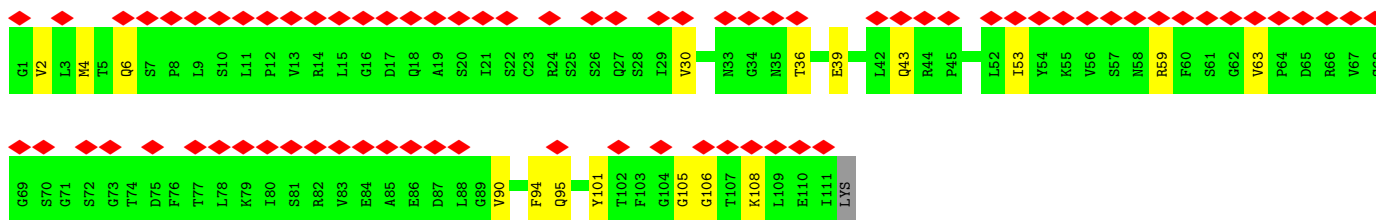
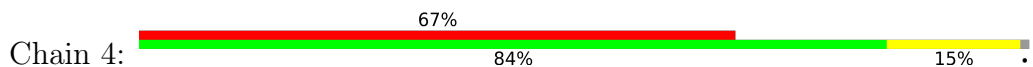
• Molecule 3: vFP7.04 Heavy chain



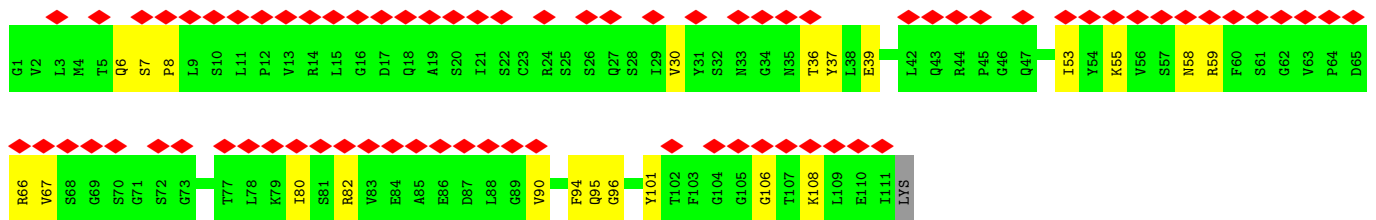
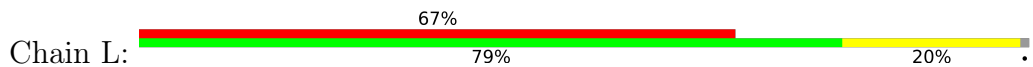
• Molecule 3: vFP7.04 Heavy chain



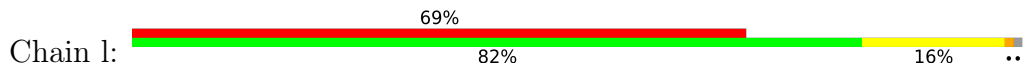
• Molecule 4: vFP7.04 light chain

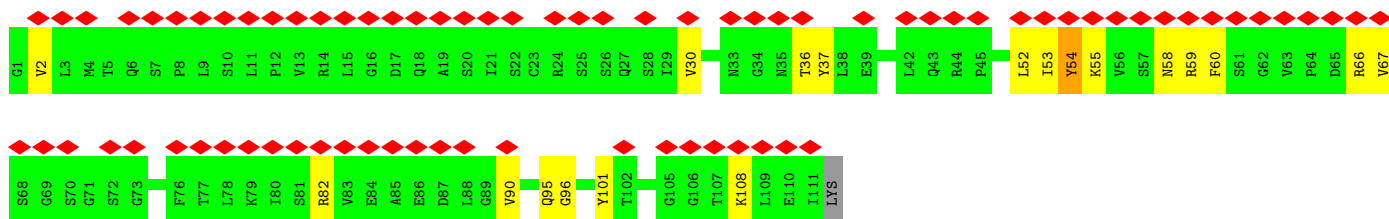


• Molecule 4: vFP7.04 light chain

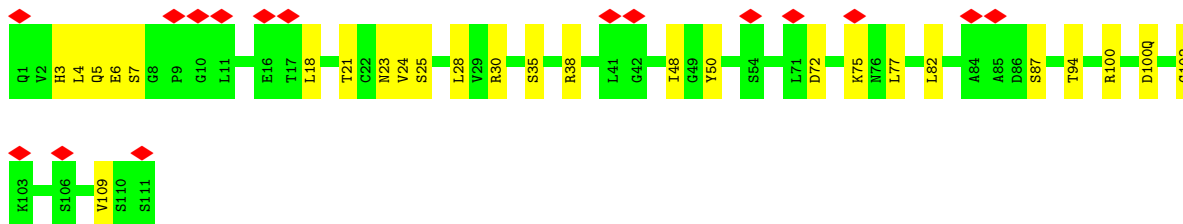
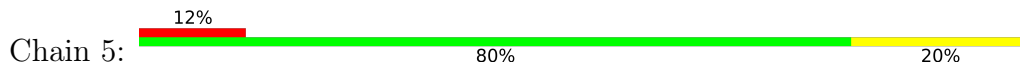


• Molecule 4: vFP7.04 light chain

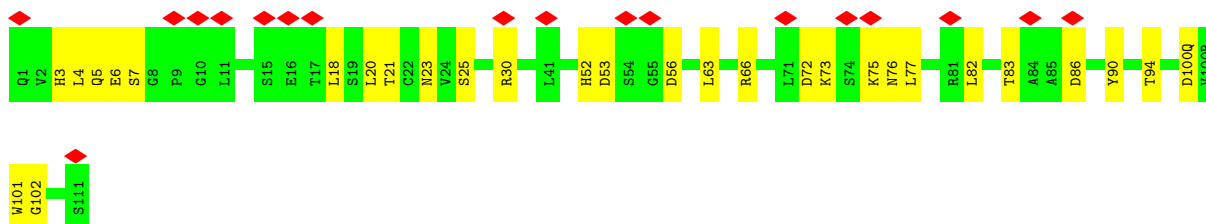
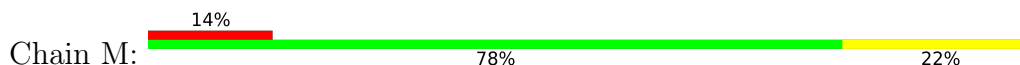




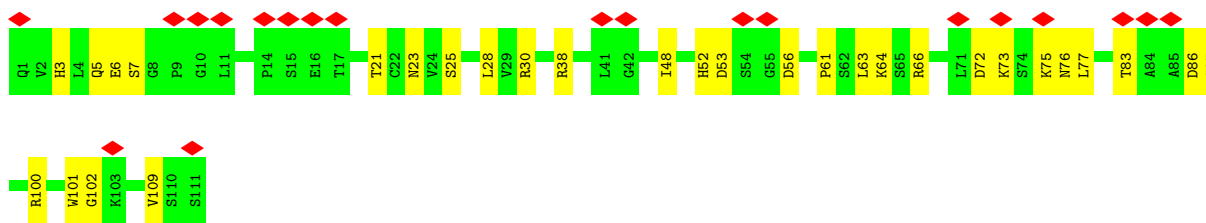
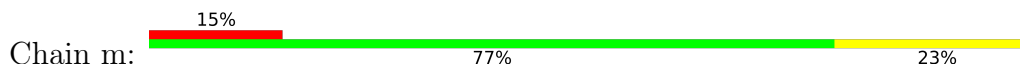
• Molecule 5: PGT122 heavy Chain



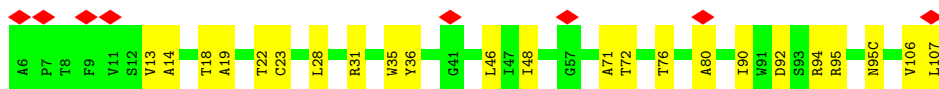
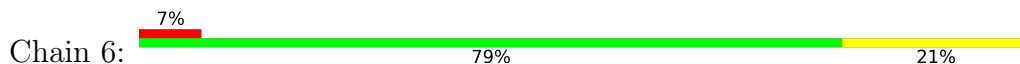
• Molecule 5: PGT122 heavy Chain



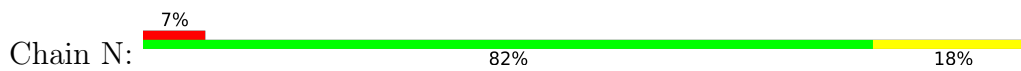
• Molecule 5: PGT122 heavy Chain



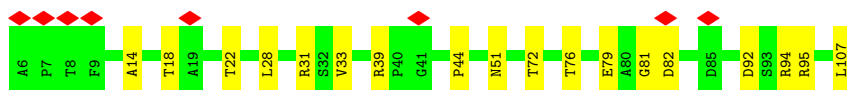
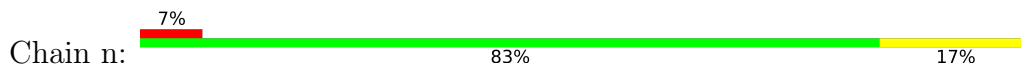
• Molecule 6: PGT122 Light chain



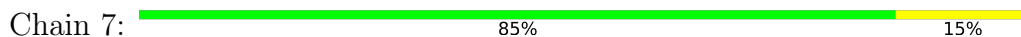
• Molecule 6: PGT122 Light chain



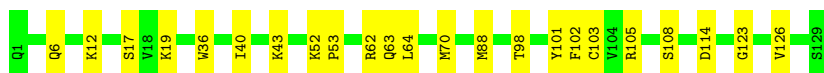
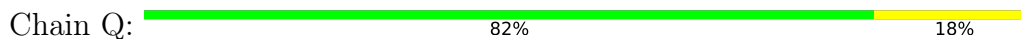
• Molecule 6: PGT122 Light chain



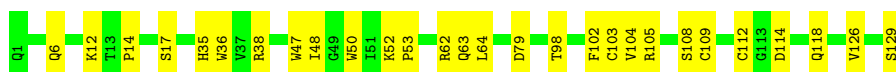
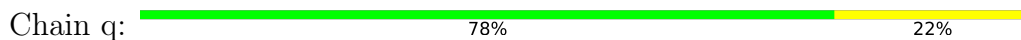
• Molecule 7: VRC03 Heavy chain



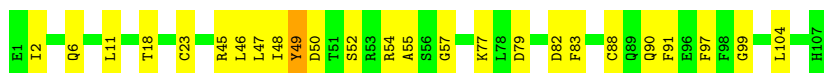
• Molecule 7: VRC03 Heavy chain



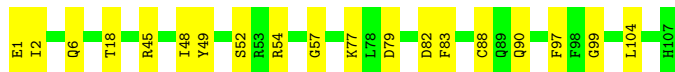
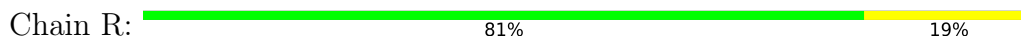
• Molecule 7: VRC03 Heavy chain




• Molecule 8: VRC03 light chain



• Molecule 8: VRC03 light chain



• Molecule 8: VRC03 light chain

Chain r:  77% 21%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain A:  50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  100%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  100%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:  100%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T: 50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain V: 50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Y: 50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z: 100%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain b: 100%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain e: 50% 50%



- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain g:  100%

MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain j:  50% 50%

MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain o:  50% 50%

MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain t:  50% 50%

MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain u:  100%

MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain w:  100%

MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain x:  100%

MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain z:  50% 50%


MAG1
MAG2

- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 9:  50% 50%

MAG1
MAG2

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain B:  100%

MAG1
MAG2
BMA3

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  100%


MAG1
MAG2
BMA3

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  67% 33%

MAG1
MAG2
BMA3

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain W:  100%

MAG1
MAG2
BMA3

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain X:  100%

MAG1
MAG2
BMA3

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain f:  67% 33%

MAG1
MAG2
BMA3

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain p:  100%

MAG1
MAG2
BMA3

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain s:  100%

MAG1
MAG2
BMA3

- Molecule 10: 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%

MAG1
MAG2
BMA3

- Molecule 11: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  40% 60%

MAG1
MAG2
BMA3
MAN4
MAN5

- Molecule 11: 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 a:  60% 40%



- Molecule 11: 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:  60% 40%

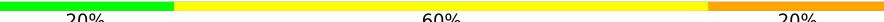


- Molecule 12: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S:  20% 60% 20%



- Molecule 12: alpha-D-mannopyranose-(1-2)-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 i:  20% 60% 20%

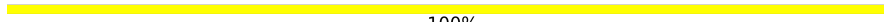


- Molecule 12: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 0:  20% 80%




- Molecule 13: 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:  100%



- Molecule 13: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain k:  75% 25%

MAG1
MAG2
BMA3
MAN4

- Molecule 13: α -D-mannopyranose-(1-3)- β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain AA:  100%

MAG1
MAG2
BMA3
MAN4

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	64580	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	53.11	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	105000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	5.517	Depositor
Minimum map value	-2.315	Depositor
Average map value	0.015	Depositor
Map value standard deviation	0.141	Depositor
Recommended contour level	1.04	Depositor
Map size (Å)	422.40002, 422.40002, 422.40002	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.1, 1.1, 1.1	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	2	0.38	0/3638	0.60	2/4939 (0.0%)
1	C	0.39	0/3638	0.60	2/4939 (0.0%)
1	c	0.38	0/3638	0.60	2/4939 (0.0%)
2	1	0.35	0/1052	0.60	0/1427
2	D	0.35	0/1052	0.59	0/1427
2	d	0.37	0/1052	0.62	0/1427
3	3	0.21	0/943	0.46	0/1282
3	H	0.21	0/943	0.46	0/1282
3	h	0.22	0/943	0.47	0/1282
4	4	0.20	0/872	0.48	0/1181
4	L	0.22	0/872	0.50	0/1181
4	l	0.20	0/872	0.54	2/1181 (0.2%)
5	5	0.22	0/1076	0.49	0/1465
5	M	0.22	0/1076	0.50	0/1465
5	m	0.22	0/1076	0.50	0/1465
6	6	0.26	0/839	0.50	0/1149
6	N	0.27	0/839	0.51	0/1149
6	n	0.25	0/839	0.49	0/1149
7	7	0.40	0/1062	0.59	0/1447
7	Q	0.42	0/1062	0.56	0/1447
7	q	0.41	0/1062	0.55	0/1447
8	8	0.36	0/820	0.69	2/1107 (0.2%)
8	R	0.35	0/820	0.71	2/1107 (0.2%)
8	r	0.34	0/820	0.70	2/1107 (0.2%)
All	All	0.33	0/30906	0.57	14/41991 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	2	0	1
1	c	0	1
8	8	0	3
8	R	0	1
8	r	0	1
All	All	0	7

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
8	r	49	TYR	CA-C-N	5.72	132.47	121.54
8	r	49	TYR	C-N-CA	5.72	132.47	121.54
1	C	99	ASN	CA-C-N	5.55	132.15	121.54
1	C	99	ASN	C-N-CA	5.55	132.15	121.54
8	R	49	TYR	CA-C-N	5.55	132.13	121.54
8	R	49	TYR	C-N-CA	5.55	132.13	121.54
1	c	99	ASN	CA-C-N	5.40	131.85	121.54
1	c	99	ASN	C-N-CA	5.40	131.85	121.54
8	8	49	TYR	CA-C-N	5.38	131.82	121.54
8	8	49	TYR	C-N-CA	5.38	131.82	121.54
4	l	54	TYR	CA-C-N	5.23	131.53	121.54
4	l	54	TYR	C-N-CA	5.23	131.53	121.54
1	2	99	ASN	CA-C-N	5.06	131.21	121.54
1	2	99	ASN	C-N-CA	5.06	131.21	121.54

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	2	77	THR	Peptide
8	8	50	ASP	Peptide
8	8	83	PHE	Peptide
8	8	90	GLN	Peptide
8	R	90	GLN	Peptide
1	c	77	THR	Peptide
8	r	90	GLN	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	2	3564	0	3489	42	0
1	C	3564	0	3489	41	0
1	c	3564	0	3489	45	0
2	1	1034	0	1014	15	0
2	D	1034	0	1014	18	0
2	d	1034	0	1014	19	0
3	3	919	0	891	17	0
3	H	919	0	891	17	0
3	h	919	0	891	17	0
4	4	853	0	841	11	0
4	L	853	0	841	14	0
4	l	853	0	841	10	0
5	5	1047	0	1026	15	0
5	M	1047	0	1026	18	0
5	m	1047	0	1026	18	0
6	6	817	0	773	13	0
6	N	817	0	773	11	0
6	n	817	0	773	11	0
7	7	1029	0	974	12	0
7	Q	1029	0	974	15	0
7	q	1029	0	974	19	0
8	8	802	0	778	14	0
8	R	802	0	778	11	0
8	r	802	0	778	15	0
9	9	28	0	25	0	0
9	A	28	0	25	0	0
9	F	28	0	25	0	0
9	G	28	0	25	0	0
9	J	28	0	25	0	0
9	K	28	0	25	0	0
9	P	28	0	25	0	0
9	T	28	0	25	0	0
9	V	28	0	25	0	0
9	Y	28	0	25	1	0
9	Z	28	0	25	0	0
9	b	28	0	25	0	0
9	e	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	g	28	0	25	0	0
9	j	28	0	25	0	0
9	o	28	0	25	0	0
9	t	28	0	25	1	0
9	u	28	0	25	0	0
9	w	28	0	25	0	0
9	x	28	0	25	0	0
9	z	28	0	25	0	0
10	B	39	0	34	1	0
10	E	39	0	34	0	0
10	O	39	0	34	0	0
10	W	39	0	34	1	0
10	X	39	0	34	0	0
10	f	39	0	34	0	0
10	p	39	0	34	1	0
10	s	39	0	34	0	0
10	y	39	0	34	0	0
11	I	61	0	52	0	0
11	a	61	0	52	0	0
11	v	61	0	52	0	0
12	0	61	0	52	0	0
12	S	61	0	52	1	0
12	i	61	0	52	1	0
13	AA	50	0	43	0	0
13	U	50	0	43	0	0
13	k	50	0	43	1	0
14	2	14	0	13	0	0
14	C	14	0	13	0	0
14	c	14	0	13	0	0
All	All	31692	0	30669	385	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (385) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:c:457:ASP:OD2	1:c:469:ARG:NH1	2.13	0.81
2:D:542:ARG:NH2	2:d:647:GLU:OE2	2.12	0.81
3:3:94:TYR:O	3:3:111:GLY:HA2	1.81	0.80
5:M:83:THR:N	5:M:86:ASP:OD2	2.18	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:h:100:ARG:HB2	3:h:103:TRP:HB2	1.69	0.74
3:H:69:MET:HB2	3:H:82:GLU:HB2	1.70	0.73
3:H:94:TYR:O	3:H:111:GLY:HA2	1.90	0.70
3:H:100:ARG:HB2	3:H:103:TRP:HB2	1.73	0.70
5:m:83:THR:N	5:m:86:ASP:OD2	2.19	0.70
2:1:647:GLU:OE2	2:d:542:ARG:NH2	2.18	0.70
2:1:542:ARG:NH2	2:D:647:GLU:OE2	2.17	0.69
3:3:38:LYS:HB2	3:3:46:GLU:HB2	1.74	0.69
3:3:20:LEU:HB2	3:3:81:MET:HB2	1.75	0.69
4:L:53:ILE:HG12	4:L:59:ARG:HG2	1.73	0.69
3:3:100:ARG:HB2	3:3:103:TRP:HB2	1.75	0.68
3:3:69:MET:HB2	3:3:82:GLU:HB2	1.76	0.68
2:d:618:ASN:N	2:d:621:GLU:OE2	2.24	0.68
2:D:618:ASN:N	2:D:621:GLU:OE2	2.26	0.68
3:h:38:LYS:HB2	3:h:46:GLU:HB2	1.76	0.68
3:H:38:LYS:HB2	3:H:46:GLU:HB2	1.76	0.67
5:m:5:GLN:HB3	5:m:23:ASN:HB2	1.77	0.67
3:h:94:TYR:O	3:h:111:GLY:HA2	1.95	0.66
4:l:37:TYR:HB2	4:l:96:GLY:HA3	1.78	0.66
7:7:109:CYS:SG	7:7:112:CYS:N	2.70	0.65
1:C:95:MET:HE1	1:C:273:ARG:HD3	1.79	0.64
8:8:18:THR:HG22	8:8:77:LYS:H	1.61	0.64
1:C:249:HIS:HD1	1:C:486:TYR:HH	1.46	0.64
3:h:10:GLU:HB2	3:h:114:VAL:HG22	1.78	0.63
7:q:109:CYS:SG	7:q:112:CYS:N	2.71	0.63
8:R:79:ASP:H	8:R:82:ASP:HB2	1.63	0.63
3:3:2:VAL:HA	3:3:25:SER:O	1.99	0.63
4:l:53:ILE:HA	4:l:59:ARG:HA	1.82	0.62
1:2:37:THR:HG22	2:1:605:CYS:HA	1.82	0.62
1:C:37:THR:HG22	2:D:605:CYS:HA	1.82	0.62
1:C:476:ARG:HA	1:C:479:TRP:HD1	1.65	0.62
1:C:373:THR:OG1	1:C:374:HIS:N	2.32	0.62
4:L:95:GLN:HE21	4:L:101:TYR:HA	1.65	0.62
3:h:69:MET:HB2	3:h:82:GLU:HB2	1.82	0.61
1:2:95:MET:HE1	1:2:273:ARG:HD3	1.82	0.61
2:1:640:GLN:HA	2:1:643:TYR:HB2	1.82	0.61
7:q:105:ARG:HH22	7:q:108:SER:HB3	1.64	0.61
1:c:491:ILE:O	2:d:585:ARG:NH2	2.34	0.61
8:8:6:GLN:HE22	8:8:99:GLY:HA3	1.66	0.61
5:M:72:ASP:HB3	5:M:77:LEU:H	1.65	0.61
5:5:3:HIS:H	5:5:25:SER:HB2	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:4:53:ILE:HA	4:4:59:ARG:HA	1.83	0.60
1:C:503:ARG:HH12	2:D:597:GLY:HA3	1.67	0.60
6:n:39:ARG:NH1	6:n:81:GLY:O	2.35	0.60
4:l:90:VAL:HG22	4:l:108:LYS:HG2	1.83	0.59
1:c:63:THR:O	1:c:67:ASN:ND2	2.35	0.59
1:2:476:ARG:HA	1:2:479:TRP:HD1	1.67	0.59
1:C:365:SER:OG	1:C:469:ARG:NH1	2.35	0.59
1:2:491:ILE:O	2:1:585:ARG:NH2	2.35	0.59
6:N:22:THR:HG22	6:N:72:THR:HG22	1.83	0.59
1:c:373:THR:OG1	1:c:374:HIS:N	2.36	0.59
3:3:18:VAL:HB	3:3:83:LEU:HB2	1.85	0.58
1:c:37:THR:HG22	2:d:605:CYS:HA	1.85	0.58
1:C:491:ILE:O	2:D:585:ARG:NH2	2.37	0.58
3:H:51:ILE:HD12	3:H:70:LEU:HB3	1.85	0.58
7:Q:63:GLN:HG2	7:Q:64:LEU:HD12	1.85	0.58
3:h:51:ILE:HD12	3:h:70:LEU:HB3	1.84	0.58
1:c:476:ARG:HA	1:c:479:TRP:HD1	1.67	0.58
3:H:39:GLN:HB3	3:H:93:VAL:HB	1.85	0.58
1:c:279:ASN:ND2	7:q:114:ASP:OD2	2.33	0.58
2:d:640:GLN:HA	2:d:643:TYR:HB2	1.86	0.57
2:1:591:GLN:NE2	2:d:541:ALA:O	2.37	0.57
5:5:5:GLN:HB3	5:5:23:ASN:HB2	1.86	0.57
8:R:48:ILE:HA	8:R:54:ARG:HA	1.86	0.57
1:C:63:THR:O	1:C:67:ASN:ND2	2.37	0.57
6:6:22:THR:HG22	6:6:72:THR:HG22	1.86	0.57
5:m:30:ARG:HH21	5:m:73:LYS:HD3	1.68	0.57
7:Q:12:LYS:NZ	7:Q:17:SER:O	2.38	0.57
1:2:258:GLN:HB2	1:2:374:HIS:HB3	1.87	0.57
3:h:93:VAL:HG22	3:h:113:THR:HG22	1.87	0.56
1:C:264:SER:O	1:C:287:GLN:NE2	2.38	0.56
2:D:541:ALA:O	2:d:591:GLN:NE2	2.38	0.56
5:M:52:HIS:HB3	5:M:56:ASP:HB3	1.87	0.56
3:3:10:GLU:HB2	3:3:114:VAL:HG22	1.87	0.56
8:8:6:GLN:NE2	8:8:88:CYS:SG	2.79	0.56
7:q:63:GLN:HG2	7:q:64:LEU:HD12	1.87	0.56
2:D:512:ALA:N	3:H:103:TRP:O	2.39	0.56
6:n:22:THR:HG22	6:n:72:THR:HG22	1.86	0.55
4:4:90:VAL:HG22	4:4:108:LYS:HG2	1.88	0.55
5:M:63:LEU:HD22	5:M:66:ARG:HE	1.71	0.55
1:2:63:THR:O	1:2:67:ASN:ND2	2.39	0.55
1:2:370:GLU:HB3	1:2:384:TYR:HE1	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:5:72:ASP:HB3	5:5:77:LEU:H	1.70	0.55
8:R:6:GLN:HE22	8:R:99:GLY:HA3	1.72	0.55
7:7:63:GLN:HG2	7:7:64:LEU:HD12	1.87	0.55
8:R:45:ARG:NH2	8:R:57:GLY:O	2.39	0.55
1:2:324:GLY:O	6:6:94:ARG:NH2	2.40	0.55
4:l:2:VAL:O	4:l:95:GLN:NE2	2.40	0.55
1:2:139:THR:HG22	10:W:1:NAG:H61	1.86	0.55
1:C:139:THR:HG22	10:p:1:NAG:H61	1.89	0.55
1:C:324:GLY:O	6:N:94:ARG:NH2	2.40	0.55
1:c:324:GLY:O	6:n:94:ARG:NH2	2.40	0.55
6:N:18:THR:HG22	6:N:76:THR:HG22	1.88	0.55
1:2:279:ASN:ND2	7:7:114:ASP:OD2	2.35	0.54
5:5:7:SER:OG	5:5:21:THR:OG1	2.25	0.54
4:L:53:ILE:HA	4:L:59:ARG:HA	1.90	0.54
1:c:139:THR:HG22	10:B:1:NAG:H61	1.88	0.54
8:R:6:GLN:NE2	8:R:88:CYS:SG	2.80	0.54
2:1:541:ALA:O	2:D:591:GLN:NE2	2.40	0.54
4:4:30:VAL:HA	4:4:36:THR:HG22	1.90	0.54
1:C:467:THR:H	7:Q:62:ARG:HH12	1.55	0.54
6:6:90:ILE:O	6:6:95(C):ASN:ND2	2.41	0.54
5:M:72:ASP:O	5:M:76:ASN:HA	2.08	0.54
5:M:72:ASP:OD2	5:M:75:LYS:HB2	2.08	0.54
5:m:7:SER:OG	5:m:21:THR:OG1	2.26	0.54
1:C:279:ASN:ND2	7:Q:114:ASP:OD2	2.31	0.53
6:N:31:ARG:HG2	6:N:92:ASP:HA	1.90	0.53
4:l:30:VAL:HA	4:l:36:THR:HG22	1.89	0.53
3:H:93:VAL:HG22	3:H:113:THR:HG22	1.90	0.53
4:L:30:VAL:HA	4:L:36:THR:HG22	1.90	0.53
6:N:14:ALA:HA	6:N:107:LEU:HB2	1.90	0.53
3:H:23:LYS:HG2	3:H:78:THR:HG22	1.91	0.53
1:2:277:ILE:O	1:2:456:ARG:NH1	2.37	0.53
1:2:350:ARG:NH2	1:2:396:ILE:O	2.40	0.53
3:H:52:VAL:HB	3:H:57:PHE:HB3	1.91	0.53
8:r:6:GLN:NE2	8:r:88:CYS:SG	2.81	0.53
4:4:59:ARG:HD2	4:4:63:VAL:HB	1.91	0.53
3:H:22:CYS:HB3	3:H:79:ALA:HB3	1.90	0.53
5:m:28:LEU:HA	5:m:76:ASN:HD21	1.73	0.53
5:m:87:SER:HB2	5:m:109:VAL:HG12	1.91	0.52
1:C:277:ILE:O	1:C:456:ARG:NH1	2.38	0.52
1:2:268:GLU:O	1:2:289:ASN:ND2	2.42	0.52
3:3:27:TYR:HB2	3:3:98:ARG:HH12	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:d:615:SER:O	2:d:617:ARG:NH1	2.43	0.52
1:c:268:GLU:O	1:c:289:ASN:ND2	2.42	0.52
4:L:67:VAL:HG22	4:L:80:ILE:HG12	1.90	0.52
8:r:45:ARG:NH2	8:r:57:GLY:O	2.43	0.52
1:c:462:ASN:N	8:r:1:GLU:OE2	2.42	0.52
1:2:304:ARG:N	1:2:440:GLN:OE1	2.42	0.52
6:6:18:THR:HG22	6:6:76:THR:HG22	1.92	0.52
6:N:39:ARG:NH1	6:N:81:GLY:O	2.36	0.52
5:m:53:ASP:OD2	5:m:73:LYS:NZ	2.40	0.52
6:n:14:ALA:HA	6:n:107:LEU:HB2	1.90	0.52
6:n:18:THR:HG22	6:n:76:THR:HG22	1.91	0.52
1:2:264:SER:O	1:2:287:GLN:NE2	2.42	0.52
1:C:44:VAL:HG23	2:D:629:LEU:HD23	1.91	0.52
6:n:33:VAL:HG12	6:n:51:ASN:HA	1.92	0.52
3:3:23:LYS:HG2	3:3:78:THR:HG22	1.92	0.52
7:q:6:GLN:HE22	7:q:102:PHE:HA	1.74	0.52
8:R:18:THR:HG22	8:R:77:LYS:H	1.73	0.51
4:l:52:LEU:HD13	4:l:67:VAL:HG11	1.92	0.51
8:r:61:ARG:NH2	8:r:82:ASP:OD1	2.39	0.51
8:R:79:ASP:HB3	8:R:82:ASP:H	1.75	0.51
5:5:72:ASP:OD2	5:5:75:LYS:HB2	2.10	0.51
8:r:6:GLN:HE22	8:r:99:GLY:HA3	1.75	0.51
1:c:95:MET:HE3	1:c:235:GLY:HA3	1.92	0.51
3:3:93:VAL:HG22	3:3:113:THR:HG22	1.92	0.51
5:m:3:HIS:H	5:m:25:SER:HB2	1.76	0.51
2:D:530:MET:HE2	2:D:623:TRP:HB3	1.91	0.51
6:n:31:ARG:HG2	6:n:92:ASP:HA	1.93	0.50
6:6:13:VAL:HG11	6:6:19:ALA:HB2	1.94	0.50
7:Q:105:ARG:HH22	7:Q:108:SER:HB3	1.75	0.50
4:4:6:GLN:OE1	4:4:106:GLY:N	2.44	0.50
8:8:45:ARG:NH2	8:8:57:GLY:O	2.45	0.50
3:H:24:ALA:HB1	3:H:98:ARG:HH21	1.76	0.50
1:2:385:CYS:HA	1:2:418:CYS:HA	1.94	0.50
5:M:7:SER:OG	5:M:21:THR:OG1	2.26	0.50
8:R:48:ILE:HD11	8:R:52:SER:HA	1.93	0.50
1:C:92:GLU:HA	1:C:238:PRO:HA	1.92	0.50
1:c:122:LEU:HD11	1:c:203:GLN:HB2	1.94	0.50
3:3:51:ILE:HD12	3:3:70:LEU:HB3	1.94	0.50
6:6:14:ALA:HA	6:6:107:LEU:HB2	1.94	0.50
2:D:640:GLN:HA	2:D:643:TYR:HB2	1.93	0.50
5:M:18:LEU:HB3	5:M:82:LEU:HB3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:M:20:LEU:HD11	5:M:90:TYR:HD2	1.77	0.50
6:N:38:GLN:NE2	6:N:42:GLN:O	2.44	0.50
3:h:20:LEU:HD12	3:h:81:MET:HG3	1.93	0.50
1:2:491:ILE:HD11	2:1:523:LEU:HD11	1.94	0.49
4:4:4:MET:SD	4:4:95:GLN:NE2	2.81	0.49
8:8:6:GLN:HG3	8:8:23:CYS:HB3	1.94	0.49
1:C:299:PRO:HA	1:C:442:VAL:HG13	1.94	0.49
3:h:39:GLN:HB3	3:h:93:VAL:HB	1.94	0.49
1:c:370:GLU:HB3	1:c:384:TYR:HE1	1.78	0.49
1:C:171:LYS:HE3	9:t:1:NAG:H81	1.94	0.49
1:c:46:LYS:HE2	1:c:492:GLU:OE2	2.12	0.49
8:r:79:ASP:H	8:r:82:ASP:HB2	1.78	0.49
1:2:373:THR:OG1	1:2:374:HIS:N	2.46	0.49
1:C:385:CYS:HA	1:C:418:CYS:HA	1.95	0.49
8:8:47:LEU:HG	8:8:48:ILE:HG22	1.95	0.49
7:Q:6:GLN:HE22	7:Q:102:PHE:HA	1.78	0.49
7:7:12:LYS:NZ	7:7:17:SER:O	2.46	0.48
1:c:385:CYS:HA	1:c:418:CYS:HA	1.95	0.48
5:5:94:THR:OG1	5:5:100(Q):ASP:OD1	2.28	0.48
2:D:615:SER:O	2:D:617:ARG:NH1	2.46	0.48
1:c:40:TYR:HB3	2:d:602:LEU:HD12	1.94	0.48
3:h:91:SER:HA	3:h:114:VAL:O	2.12	0.48
5:m:63:LEU:HD22	5:m:66:ARG:HE	1.79	0.48
7:q:12:LYS:NZ	7:q:17:SER:O	2.47	0.48
1:c:137:ASN:N	6:n:94:ARG:O	2.46	0.48
7:7:6:GLN:HE22	7:7:102:PHE:HA	1.78	0.48
2:d:638:TYR:HB2	2:d:641:ILE:HD12	1.95	0.48
1:C:270:VAL:HG11	1:C:345:VAL:HG12	1.93	0.48
8:r:18:THR:HG22	8:r:77:LYS:H	1.78	0.48
1:c:272:ILE:HG22	1:c:286:VAL:HG22	1.95	0.48
3:H:20:LEU:HD12	3:H:81:MET:HG3	1.96	0.48
6:n:28:LEU:HD13	6:n:95:ARG:HE	1.79	0.48
8:r:37:GLN:HB2	8:r:47:LEU:HD22	1.95	0.48
5:5:87:SER:HB2	5:5:109:VAL:HG12	1.95	0.47
7:Q:98:THR:HA	7:Q:126:VAL:O	2.14	0.47
5:5:38:ARG:HB3	5:5:48:ILE:HD11	1.97	0.47
8:8:46:LEU:HD21	8:8:49:TYR:HB3	1.95	0.47
8:8:48:ILE:HD11	8:8:52:SER:HA	1.95	0.47
1:C:58:ALA:HA	1:C:61:TYR:HD2	1.79	0.47
3:3:48:ILE:HG21	3:3:81:MET:HE1	1.96	0.47
3:H:91:SER:HA	3:H:114:VAL:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:37:TYR:HB2	4:L:96:GLY:HA3	1.97	0.47
7:7:36:TRP:HE1	7:7:103:CYS:HB3	1.79	0.47
1:2:460:SER:OG	7:7:63:GLN:NE2	2.48	0.47
1:c:304:ARG:N	1:c:440:GLN:OE1	2.46	0.47
6:6:28:LEU:HD13	6:6:95:ARG:HE	1.80	0.47
2:d:620:SER:O	2:d:624:ASP:N	2.48	0.47
7:q:118:GLN:HE22	8:r:55:ALA:HB1	1.80	0.47
5:5:6:GLU:OE2	5:5:102:GLY:HA3	2.15	0.47
5:m:72:ASP:OD2	5:m:75:LYS:HB2	2.15	0.47
1:c:95:MET:HE1	1:c:273:ARG:HD3	1.96	0.46
1:2:137:ASN:N	6:6:94:ARG:O	2.48	0.46
1:2:249:HIS:ND1	1:2:486:TYR:OH	2.41	0.46
7:Q:19:LYS:HA	7:Q:88:MET:O	2.16	0.46
7:Q:36:TRP:HE1	7:Q:103:CYS:HB3	1.80	0.46
3:h:52:VAL:HB	3:h:57:PHE:HB3	1.97	0.46
1:C:272:ILE:HG22	1:C:286:VAL:HG22	1.97	0.46
5:M:30:ARG:HH21	5:M:73:LYS:HD3	1.79	0.46
1:c:460:SER:OG	7:q:63:GLN:NE2	2.44	0.46
1:2:342:LEU:HA	1:2:345:VAL:HG22	1.97	0.46
1:C:312:GLY:H	1:C:315:GLN:HB2	1.80	0.46
7:q:98:THR:HA	7:q:126:VAL:O	2.14	0.46
8:r:15:PRO:HD3	8:r:106:VAL:HG13	1.98	0.46
3:3:91:SER:HA	3:3:114:VAL:O	2.16	0.46
7:7:118:GLN:HE22	8:8:55:ALA:HB1	1.81	0.46
5:m:6:GLU:OE2	5:m:102:GLY:HA3	2.15	0.46
8:r:48:ILE:HD11	8:r:52:SER:HA	1.97	0.46
1:c:31:ALA:N	1:c:33:ASN:OD1	2.49	0.46
8:8:11:LEU:HD23	8:8:104:LEU:HG	1.98	0.46
5:m:38:ARG:HB3	5:m:48:ILE:HD11	1.97	0.46
2:D:533:ALA:HA	2:D:536:THR:HG22	1.98	0.46
5:m:28:LEU:HG	5:m:30:ARG:H	1.80	0.46
1:2:92:GLU:HA	1:2:238:PRO:HA	1.97	0.46
3:3:24:ALA:HB3	3:3:29:PHE:HE1	1.81	0.46
8:r:46:LEU:HD21	8:r:49:TYR:HB3	1.97	0.45
1:2:49:GLU:HB3	1:2:99:ASN:HD22	1.81	0.45
1:C:40:TYR:HB3	2:D:602:LEU:HD12	1.97	0.45
1:C:259:LEU:HD23	1:C:449:ILE:HD13	1.98	0.45
3:H:24:ALA:HB3	3:H:29:PHE:HE1	1.82	0.45
6:6:31:ARG:HG2	6:6:92:ASP:HA	1.98	0.45
1:2:100:MET:HG3	1:2:483:LEU:HD13	1.99	0.45
8:R:83:PHE:HE1	8:R:104:LEU:H	1.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:312:GLY:H	1:2:315:GLN:HB2	1.81	0.45
2:d:577:GLN:HA	2:d:580:VAL:HG22	1.97	0.45
1:c:299:PRO:HA	1:c:442:VAL:HG13	1.99	0.45
1:c:463:SER:OG	7:q:62:ARG:NE	2.50	0.45
1:C:491:ILE:HD11	2:D:523:LEU:HD11	1.99	0.45
3:h:24:ALA:HB3	3:h:29:PHE:HE1	1.82	0.45
8:r:47:LEU:HG	8:r:48:ILE:HG22	1.99	0.45
1:C:300:ASN:ND2	1:C:327:ARG:O	2.45	0.45
4:l:54:TYR:HE1	4:l:60:PHE:HA	1.82	0.45
6:6:80:ALA:HA	6:6:106:VAL:HG11	1.99	0.45
4:L:6:GLN:OE1	4:L:106:GLY:N	2.45	0.45
5:M:6:GLU:OE2	5:M:102:GLY:HA3	2.17	0.45
1:2:281:ALA:HB3	7:7:114:ASP:OD2	2.17	0.45
8:8:91:PHE:HE1	13:k:1:NAG:H82	1.82	0.45
3:h:23:LYS:HG2	3:h:78:THR:HG22	1.98	0.45
1:c:498:PRO:HB3	2:d:610:TRP:CD2	2.52	0.44
1:2:171:LYS:HE3	9:Y:1:NAG:H81	1.99	0.44
1:C:491:ILE:HG22	1:C:493:PRO:HD3	1.98	0.44
2:D:530:MET:O	2:D:534:SER:N	2.50	0.44
5:M:5:GLN:HB3	5:M:23:ASN:HB2	1.99	0.44
8:r:2:ILE:HB	8:r:97:PHE:CD2	2.52	0.44
4:4:39:GLU:HB2	4:4:94:PHE:HB3	1.99	0.44
4:l:55:LYS:HB2	4:l:58:ASN:HB2	1.99	0.44
2:1:577:GLN:HA	2:1:580:VAL:HG22	1.98	0.44
2:1:638:TYR:HB2	2:1:641:ILE:HD12	2.00	0.44
4:L:39:GLU:HB2	4:L:94:PHE:HB3	1.99	0.44
1:C:268:GLU:O	1:C:289:ASN:ND2	2.50	0.44
1:c:174:SER:OG	1:c:175:LEU:N	2.50	0.44
7:q:52:LYS:HA	7:q:53:PRO:HD3	1.86	0.44
1:c:93:PHE:CE2	1:c:228:CYS:HB2	2.53	0.44
1:2:44:VAL:HG23	2:1:629:LEU:HD23	2.00	0.44
1:2:272:ILE:HG22	1:2:286:VAL:HG22	2.00	0.44
1:C:501:CYS:HB2	2:d:661:LEU:HB3	2.00	0.44
7:Q:40:ILE:HB	7:Q:43:LYS:HB2	2.00	0.44
1:c:44:VAL:HG23	2:d:629:LEU:HD23	2.00	0.44
1:c:305:LYS:HE3	1:c:321:GLY:HA2	2.00	0.44
1:c:342:LEU:HA	1:c:345:VAL:HG22	2.00	0.43
5:5:100:ARG:NH2	12:i:4:MAN:O6	2.44	0.43
5:M:52:HIS:ND1	5:M:53:ASP:O	2.49	0.43
3:h:22:CYS:HB3	3:h:79:ALA:HB3	1.99	0.43
1:C:460:SER:OG	7:Q:63:GLN:NE2	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1:580:VAL:HA	2:1:583:VAL:HG22	2.00	0.43
7:7:35:HIS:HB2	7:7:104:VAL:HG23	1.99	0.43
3:h:47:TRP:CZ2	4:l:101:TYR:HB2	2.54	0.43
1:c:270:VAL:HG11	1:c:345:VAL:HG12	2.00	0.43
5:M:101:TRP:CE2	6:N:44:PRO:HB2	2.53	0.43
5:m:21:THR:HB	5:m:77:LEU:HD23	2.01	0.43
7:q:14:PRO:HD3	7:q:129:SER:HB3	2.00	0.43
1:c:184:ILE:HD12	1:c:184:ILE:HA	1.94	0.43
1:c:459:GLY:HA3	7:q:62:ARG:HB2	2.01	0.43
1:2:423:ILE:HG22	1:2:435:TYR:HA	2.00	0.43
3:3:30:THR:HB	3:3:54:GLU:OE2	2.18	0.43
1:c:281:ALA:HB3	7:q:114:ASP:OD2	2.18	0.43
1:c:312:GLY:H	1:c:315:GLN:HB2	1.84	0.43
1:2:122:LEU:HD11	1:2:203:GLN:HB2	2.01	0.43
1:2:259:LEU:HD23	1:2:449:ILE:HD13	2.00	0.43
1:C:164:GLU:OE2	1:C:308:ARG:HD3	2.17	0.43
1:C:174:SER:OG	1:C:175:LEU:N	2.51	0.43
5:5:18:LEU:HB3	5:5:82:LEU:HB3	2.01	0.43
5:m:52:HIS:HB3	5:m:56:ASP:HB3	2.01	0.43
5:m:101:TRP:CE2	6:n:44:PRO:HB2	2.54	0.43
7:q:35:HIS:HB2	7:q:104:VAL:HG23	2.00	0.43
1:c:264:SER:O	1:c:287:GLN:NE2	2.51	0.43
5:5:35:SER:HA	5:5:50:TYR:HA	2.01	0.43
5:m:61:PRO:HA	5:m:64:LYS:HB2	2.00	0.43
4:4:2:VAL:O	4:4:95:GLN:NE2	2.50	0.43
2:D:530:MET:HG2	2:D:623:TRP:HE3	1.84	0.43
2:d:533:ALA:HA	2:d:536:THR:HG22	2.01	0.43
7:q:47:TRP:HZ2	7:q:50:TRP:HD1	1.66	0.43
1:C:318:TYR:OH	7:q:79:ASP:OD2	2.24	0.43
1:c:61:TYR:OH	1:c:74:CYS:O	2.37	0.42
6:6:36:TYR:HE1	6:6:46:LEU:HD12	1.84	0.42
7:7:47:TRP:HZ2	7:7:50:TRP:HD1	1.66	0.42
6:n:79:GLU:HG2	6:n:82:ASP:OD2	2.19	0.42
4:4:43:GLN:HB3	4:4:90:VAL:HB	2.01	0.42
5:5:28:LEU:HG	5:5:30:ARG:H	1.83	0.42
8:8:79:ASP:H	8:8:82:ASP:HB2	1.83	0.42
8:8:2:ILE:HB	8:8:97:PHE:CD2	2.55	0.42
6:N:80:ALA:HA	6:N:106:VAL:HG11	2.02	0.42
7:Q:36:TRP:HZ3	7:Q:70:MET:HG2	1.83	0.42
1:2:67:ASN:OD1	1:2:67:ASN:N	2.53	0.42
1:2:270:VAL:HG11	1:2:345:VAL:HG12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:93:PHE:CE2	1:C:228:CYS:HB2	2.54	0.42
3:h:51:ILE:HG13	3:h:58:THR:HG22	2.02	0.42
1:2:39:TYR:O	1:2:494:LEU:HA	2.20	0.42
1:2:164:GLU:OE2	1:2:308:ARG:HD3	2.20	0.42
8:r:90:GLN:O	8:r:97:PHE:N	2.52	0.42
1:c:378:CYS:HB3	1:c:383:PHE:HE2	1.85	0.42
1:c:503:ARG:HH12	2:d:597:GLY:HA3	1.85	0.42
1:C:462:ASN:N	8:R:1:GLU:OE2	2.53	0.42
1:2:40:TYR:HB3	2:1:602:LEU:HD12	2.02	0.42
3:3:47:TRP:CZ2	4:4:101:TYR:HB2	2.55	0.42
7:Q:52:LYS:HA	7:Q:53:PRO:HD3	1.93	0.42
1:c:82:GLN:HG2	1:c:246:GLN:HB2	2.03	0.41
3:H:36:TRP:HB3	3:H:48:ILE:HD12	2.02	0.41
7:q:36:TRP:HE1	7:q:103:CYS:HB3	1.85	0.41
3:H:47:TRP:CZ2	4:L:101:TYR:HB2	2.55	0.41
4:L:90:VAL:HG22	4:L:108:LYS:HG2	2.02	0.41
6:N:60:ASP:OD1	6:N:60:ASP:N	2.53	0.41
5:m:100:ARG:NH2	12:S:4:MAN:O6	2.45	0.41
1:c:350:ARG:NH2	1:c:396:ILE:O	2.40	0.41
5:5:23:ASN:HA	5:5:77:LEU:HG	2.02	0.41
4:L:55:LYS:HD2	4:L:58:ASN:HB2	2.03	0.41
7:q:38:ARG:HB3	7:q:48:ILE:HD11	2.01	0.41
8:8:48:ILE:HA	8:8:54:ARG:HA	2.02	0.41
5:M:94:THR:OG1	5:M:100(Q):ASP:OD1	2.38	0.41
6:N:33:VAL:HG12	6:N:51:ASN:HA	2.03	0.41
1:c:274:SER:OG	1:c:275:GLU:N	2.53	0.41
7:7:98:THR:HA	7:7:126:VAL:O	2.20	0.41
1:2:437:PRO:HA	1:2:438:PRO:HD3	1.96	0.41
1:c:491:ILE:HD11	2:d:523:LEU:HD11	2.03	0.41
1:2:377:ASN:OD1	1:2:381:GLU:N	2.54	0.41
6:6:23:CYS:O	6:6:71:ALA:HB3	2.20	0.41
8:R:2:ILE:HB	8:R:97:PHE:CD2	2.56	0.41
1:C:373:THR:OG1	1:C:385:CYS:N	2.54	0.41
5:5:4:LEU:HD23	5:5:24:VAL:HA	2.02	0.41
4:L:7:SER:HA	4:L:8:PRO:HA	1.84	0.41
4:l:66:ARG:HD3	4:l:82:ARG:HB2	2.03	0.41
1:C:258:GLN:HB2	1:C:374:HIS:HB3	2.03	0.40
2:1:533:ALA:HA	2:1:536:THR:HG22	2.03	0.40
4:4:6:GLN:CD	4:4:105:GLY:H	2.29	0.40
4:L:55:LYS:HD2	4:L:58:ASN:HD22	1.86	0.40
5:M:4:LEU:HA	5:M:23:ASN:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:h:30:THR:HB	3:h:54:GLU:OE2	2.21	0.40
1:c:164:GLU:OE2	1:c:308:ARG:HD3	2.20	0.40
6:6:35:TRP:HB2	6:6:48:ILE:HB	2.03	0.40
1:C:122:LEU:HD11	1:C:203:GLN:HB2	2.04	0.40
1:C:378:CYS:HB3	1:C:383:PHE:HE2	1.85	0.40
5:M:3:HIS:H	5:M:25:SER:HB2	1.86	0.40
7:Q:101:TYR:O	7:Q:123:GLY:HA2	2.22	0.40
7:Q:114:ASP:OD1	7:Q:114:ASP:N	2.55	0.40
1:2:166:ARG:HD3	1:2:166:ARG:HA	1.89	0.40
1:2:494:LEU:HD13	2:1:593:LEU:HD11	2.04	0.40
4:L:66:ARG:HD3	4:L:82:ARG:HB2	2.04	0.40
5:M:21:THR:HB	5:M:77:LEU:HD23	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	2	447/473 (94%)	410 (92%)	37 (8%)	0	100	100
1	C	447/473 (94%)	412 (92%)	35 (8%)	0	100	100
1	c	447/473 (94%)	414 (93%)	33 (7%)	0	100	100
2	1	128/153 (84%)	120 (94%)	8 (6%)	0	100	100
2	D	128/153 (84%)	121 (94%)	7 (6%)	0	100	100
2	d	128/153 (84%)	121 (94%)	7 (6%)	0	100	100
3	3	115/118 (98%)	102 (89%)	13 (11%)	0	100	100
3	H	115/118 (98%)	108 (94%)	7 (6%)	0	100	100
3	h	115/118 (98%)	107 (93%)	8 (7%)	0	100	100
4	4	109/112 (97%)	100 (92%)	9 (8%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	L	109/112 (97%)	100 (92%)	9 (8%)	0	100	100
4	l	109/112 (97%)	98 (90%)	11 (10%)	0	100	100
5	5	130/132 (98%)	118 (91%)	12 (9%)	0	100	100
5	M	130/132 (98%)	117 (90%)	13 (10%)	0	100	100
5	m	130/132 (98%)	120 (92%)	10 (8%)	0	100	100
6	6	105/107 (98%)	96 (91%)	9 (9%)	0	100	100
6	N	105/107 (98%)	97 (92%)	8 (8%)	0	100	100
6	n	105/107 (98%)	95 (90%)	10 (10%)	0	100	100
7	7	127/129 (98%)	121 (95%)	6 (5%)	0	100	100
7	Q	127/129 (98%)	121 (95%)	6 (5%)	0	100	100
7	q	127/129 (98%)	120 (94%)	7 (6%)	0	100	100
8	8	100/102 (98%)	93 (93%)	7 (7%)	0	100	100
8	R	100/102 (98%)	95 (95%)	5 (5%)	0	100	100
8	r	100/102 (98%)	96 (96%)	4 (4%)	0	100	100
All	All	3783/3978 (95%)	3502 (93%)	281 (7%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	2	405/422 (96%)	403 (100%)	2 (0%)	81	82
1	C	405/422 (96%)	403 (100%)	2 (0%)	81	82
1	c	405/422 (96%)	403 (100%)	2 (0%)	81	82
2	1	110/129 (85%)	110 (100%)	0	100	100
2	D	110/129 (85%)	110 (100%)	0	100	100
2	d	110/129 (85%)	110 (100%)	0	100	100
3	3	99/100 (99%)	99 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	H	99/100 (99%)	99 (100%)	0	100	100
3	h	99/100 (99%)	99 (100%)	0	100	100
4	4	96/97 (99%)	96 (100%)	0	100	100
4	L	96/97 (99%)	96 (100%)	0	100	100
4	l	96/97 (99%)	96 (100%)	0	100	100
5	5	116/116 (100%)	116 (100%)	0	100	100
5	M	116/116 (100%)	116 (100%)	0	100	100
5	m	116/116 (100%)	116 (100%)	0	100	100
6	6	89/89 (100%)	89 (100%)	0	100	100
6	N	89/89 (100%)	89 (100%)	0	100	100
6	n	89/89 (100%)	89 (100%)	0	100	100
7	7	109/110 (99%)	108 (99%)	1 (1%)	70	76
7	Q	109/110 (99%)	109 (100%)	0	100	100
7	q	109/110 (99%)	109 (100%)	0	100	100
8	8	86/86 (100%)	86 (100%)	0	100	100
8	R	86/86 (100%)	86 (100%)	0	100	100
8	r	86/86 (100%)	86 (100%)	0	100	100
All	All	3330/3447 (97%)	3323 (100%)	7 (0%)	85	87

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	c	255	VAL
1	c	416	LEU
1	2	373	THR
1	2	416	LEU
1	C	373	THR
1	C	416	LEU
7	7	127	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	c	289	ASN
1	c	315	GLN

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Mol	Chain	Res	Type
1	c	432	GLN
1	2	85	HIS
1	2	203	GLN
1	2	315	GLN
1	2	432	GLN
1	C	183	GLN
1	C	203	GLN
1	C	315	GLN
1	C	432	GLN
2	1	591	GLN
2	1	616	ASN
3	3	43	HIS
6	6	42	GLN
7	7	6	GLN
7	7	118	GLN
2	D	591	GLN
2	D	607	ASN
2	D	616	ASN
2	D	650	GLN
3	H	43	HIS
4	L	58	ASN
4	L	95	GLN
6	N	89	HIS
7	Q	6	GLN
7	Q	118	GLN
8	R	90	GLN
2	d	591	GLN
2	d	616	ASN
3	h	43	HIS
4	l	58	ASN
5	m	68	HIS
5	m	76	ASN
7	q	6	GLN
7	q	118	GLN
8	r	31	ASN
8	r	90	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

111 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
12	NAG	0	1	12,1	14,14,15	0.51	0	17,19,21	0.58	0
12	NAG	0	2	12	14,14,15	0.28	0	17,19,21	1.14	1 (5%)
12	BMA	0	3	12	11,11,12	1.36	2 (18%)	15,15,17	1.22	2 (13%)
12	MAN	0	4	12	11,11,12	0.84	1 (9%)	15,15,17	1.54	2 (13%)
12	MAN	0	5	12	11,11,12	0.67	0	15,15,17	1.00	2 (13%)
9	NAG	9	1	9,1	14,14,15	0.35	0	17,19,21	1.03	1 (5%)
9	NAG	9	2	9	14,14,15	0.27	0	17,19,21	0.53	0
9	NAG	A	1	9,1	14,14,15	0.57	0	17,19,21	1.03	1 (5%)
9	NAG	A	2	9	14,14,15	0.37	0	17,19,21	0.57	0
13	NAG	AA	1	1,13	14,14,15	0.25	0	17,19,21	0.80	1 (5%)
13	NAG	AA	2	13	14,14,15	0.44	0	17,19,21	1.00	2 (11%)
13	BMA	AA	3	13	11,11,12	1.16	1 (9%)	15,15,17	0.88	0
13	MAN	AA	4	13	11,11,12	0.91	1 (9%)	15,15,17	1.39	2 (13%)
10	NAG	B	1	1,10	14,14,15	0.34	0	17,19,21	0.62	0
10	NAG	B	2	10	14,14,15	0.41	0	17,19,21	0.86	1 (5%)
10	BMA	B	3	10	11,11,12	0.89	1 (9%)	15,15,17	1.80	4 (26%)
10	NAG	E	1	1,10	14,14,15	0.36	0	17,19,21	0.53	0
10	NAG	E	2	10	14,14,15	0.27	0	17,19,21	0.48	0
10	BMA	E	3	10	11,11,12	0.69	0	15,15,17	0.78	0
9	NAG	F	1	9,1	14,14,15	0.60	1 (7%)	17,19,21	0.97	1 (5%)
9	NAG	F	2	9	14,14,15	0.23	0	17,19,21	0.49	0
9	NAG	G	1	9,1	14,14,15	0.42	0	17,19,21	0.49	0
9	NAG	G	2	9	14,14,15	0.30	0	17,19,21	0.54	0
11	NAG	I	1	11,1	14,14,15	0.25	0	17,19,21	0.65	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	NAG	I	2	11	14,14,15	0.23	0	17,19,21	0.48	0
11	BMA	I	3	11	11,11,12	0.79	0	15,15,17	0.86	0
11	MAN	I	4	11	11,11,12	0.75	0	15,15,17	1.08	2 (13%)
11	MAN	I	5	11	11,11,12	0.83	0	15,15,17	1.01	2 (13%)
9	NAG	J	1	9,1	14,14,15	0.26	0	17,19,21	0.56	0
9	NAG	J	2	9	14,14,15	0.50	0	17,19,21	0.62	1 (5%)
9	NAG	K	1	9,1	14,14,15	0.19	0	17,19,21	0.54	0
9	NAG	K	2	9	14,14,15	0.24	0	17,19,21	0.53	0
10	NAG	O	1	1,10	14,14,15	0.53	0	17,19,21	0.59	0
10	NAG	O	2	10	14,14,15	0.24	0	17,19,21	0.48	0
10	BMA	O	3	10	11,11,12	1.03	1 (9%)	15,15,17	1.17	2 (13%)
9	NAG	P	1	9,1	14,14,15	0.30	0	17,19,21	0.58	0
9	NAG	P	2	9	14,14,15	0.37	0	17,19,21	0.46	0
12	NAG	S	1	12,1	14,14,15	0.56	0	17,19,21	0.61	0
12	NAG	S	2	12	14,14,15	0.26	0	17,19,21	1.16	2 (11%)
12	BMA	S	3	12	11,11,12	1.35	3 (27%)	15,15,17	1.15	1 (6%)
12	MAN	S	4	12	11,11,12	0.79	1 (9%)	15,15,17	1.39	2 (13%)
12	MAN	S	5	12	11,11,12	0.66	0	15,15,17	1.03	2 (13%)
9	NAG	T	1	9,1	14,14,15	0.25	0	17,19,21	1.06	1 (5%)
9	NAG	T	2	9	14,14,15	0.33	0	17,19,21	0.55	0
13	NAG	U	1	1,13	14,14,15	0.36	0	17,19,21	0.84	1 (5%)
13	NAG	U	2	13	14,14,15	0.43	0	17,19,21	1.12	2 (11%)
13	BMA	U	3	13	11,11,12	1.23	1 (9%)	15,15,17	0.85	0
13	MAN	U	4	13	11,11,12	0.93	1 (9%)	15,15,17	1.36	2 (13%)
9	NAG	V	1	9,1	14,14,15	0.46	0	17,19,21	1.06	1 (5%)
9	NAG	V	2	9	14,14,15	0.33	0	17,19,21	0.60	0
10	NAG	W	1	1,10	14,14,15	0.27	0	17,19,21	0.54	0
10	NAG	W	2	10	14,14,15	0.50	0	17,19,21	0.97	1 (5%)
10	BMA	W	3	10	11,11,12	0.70	0	15,15,17	1.58	2 (13%)
10	NAG	X	1	1,10	14,14,15	0.35	0	17,19,21	0.53	0
10	NAG	X	2	10	14,14,15	0.24	0	17,19,21	0.44	0
10	BMA	X	3	10	11,11,12	0.76	0	15,15,17	0.75	0
9	NAG	Y	1	9,1	14,14,15	0.71	1 (7%)	17,19,21	1.02	1 (5%)
9	NAG	Y	2	9	14,14,15	0.25	0	17,19,21	0.45	0
9	NAG	Z	1	9,1	14,14,15	0.30	0	17,19,21	0.44	0
9	NAG	Z	2	9	14,14,15	0.34	0	17,19,21	0.57	0
11	NAG	a	1	11,1	14,14,15	0.26	0	17,19,21	0.57	0
11	NAG	a	2	11	14,14,15	0.26	0	17,19,21	0.50	0
11	BMA	a	3	11	11,11,12	0.66	0	15,15,17	0.84	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	MAN	a	4	11	11,11,12	0.78	1 (9%)	15,15,17	1.16	2 (13%)
11	MAN	a	5	11	11,11,12	0.78	0	15,15,17	1.04	2 (13%)
9	NAG	b	1	9,1	14,14,15	0.21	0	17,19,21	0.60	0
9	NAG	b	2	9	14,14,15	0.46	0	17,19,21	0.57	0
9	NAG	e	1	9,1	14,14,15	0.18	0	17,19,21	0.63	1 (5%)
9	NAG	e	2	9	14,14,15	0.26	0	17,19,21	0.49	0
10	NAG	f	1	1,10	14,14,15	0.51	0	17,19,21	0.57	0
10	NAG	f	2	10	14,14,15	0.30	0	17,19,21	0.53	0
10	BMA	f	3	10	11,11,12	0.95	1 (9%)	15,15,17	1.18	2 (13%)
9	NAG	g	1	9,1	14,14,15	0.32	0	17,19,21	0.55	0
9	NAG	g	2	9	14,14,15	0.32	0	17,19,21	0.48	0
12	NAG	i	1	12,1	14,14,15	0.57	0	17,19,21	0.61	0
12	NAG	i	2	12	14,14,15	0.28	0	17,19,21	1.17	2 (11%)
12	BMA	i	3	12	11,11,12	1.30	1 (9%)	15,15,17	1.19	1 (6%)
12	MAN	i	4	12	11,11,12	0.82	1 (9%)	15,15,17	1.46	2 (13%)
12	MAN	i	5	12	11,11,12	0.73	0	15,15,17	0.97	2 (13%)
9	NAG	j	1	9,1	14,14,15	0.27	0	17,19,21	1.02	1 (5%)
9	NAG	j	2	9	14,14,15	0.29	0	17,19,21	0.57	0
13	NAG	k	1	1,13	14,14,15	0.25	0	17,19,21	0.80	1 (5%)
13	NAG	k	2	13	14,14,15	0.39	0	17,19,21	0.98	1 (5%)
13	BMA	k	3	13	11,11,12	1.06	1 (9%)	15,15,17	0.76	0
13	MAN	k	4	13	11,11,12	0.79	1 (9%)	15,15,17	1.21	2 (13%)
9	NAG	o	1	9,1	14,14,15	0.39	0	17,19,21	1.06	1 (5%)
9	NAG	o	2	9	14,14,15	0.35	0	17,19,21	0.56	0
10	NAG	p	1	1,10	14,14,15	0.25	0	17,19,21	0.57	0
10	NAG	p	2	10	14,14,15	0.46	0	17,19,21	0.96	1 (5%)
10	BMA	p	3	10	11,11,12	0.68	0	15,15,17	1.49	1 (6%)
10	NAG	s	1	1,10	14,14,15	0.39	0	17,19,21	0.54	0
10	NAG	s	2	10	14,14,15	0.26	0	17,19,21	0.48	0
10	BMA	s	3	10	11,11,12	0.72	0	15,15,17	0.76	0
9	NAG	t	1	9,1	14,14,15	0.60	1 (7%)	17,19,21	1.04	1 (5%)
9	NAG	t	2	9	14,14,15	0.28	0	17,19,21	0.43	0
9	NAG	u	1	9,1	14,14,15	0.23	0	17,19,21	0.47	0
9	NAG	u	2	9	14,14,15	0.33	0	17,19,21	0.55	0
11	NAG	v	1	11,1	14,14,15	0.29	0	17,19,21	0.57	0
11	NAG	v	2	11	14,14,15	0.23	0	17,19,21	0.46	0
11	BMA	v	3	11	11,11,12	0.73	0	15,15,17	0.85	0
11	MAN	v	4	11	11,11,12	0.87	1 (9%)	15,15,17	1.25	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	MAN	v	5	11	11,11,12	0.86	0	15,15,17	0.90	1 (6%)
9	NAG	w	1	9,1	14,14,15	0.25	0	17,19,21	0.52	0
9	NAG	w	2	9	14,14,15	0.57	0	17,19,21	0.58	0
9	NAG	x	1	9,1	14,14,15	0.18	0	17,19,21	0.57	0
9	NAG	x	2	9	14,14,15	0.27	0	17,19,21	0.48	0
10	NAG	y	1	1,10	14,14,15	0.39	0	17,19,21	0.55	0
10	NAG	y	2	10	14,14,15	0.27	0	17,19,21	0.52	0
10	BMA	y	3	10	11,11,12	1.09	1 (9%)	15,15,17	1.16	1 (6%)
9	NAG	z	1	9,1	14,14,15	0.41	0	17,19,21	0.73	1 (5%)
9	NAG	z	2	9	14,14,15	0.31	0	17,19,21	0.50	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
12	NAG	0	1	12,1	-	0/6/23/26	0/1/1/1
12	NAG	0	2	12	-	4/6/23/26	0/1/1/1
12	BMA	0	3	12	-	0/2/19/22	0/1/1/1
12	MAN	0	4	12	-	0/2/19/22	0/1/1/1
12	MAN	0	5	12	-	1/2/19/22	0/1/1/1
9	NAG	9	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	9	2	9	-	0/6/23/26	0/1/1/1
9	NAG	A	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	A	2	9	-	2/6/23/26	0/1/1/1
13	NAG	AA	1	1,13	-	2/6/23/26	0/1/1/1
13	NAG	AA	2	13	-	2/6/23/26	0/1/1/1
13	BMA	AA	3	13	-	0/2/19/22	0/1/1/1
13	MAN	AA	4	13	-	2/2/19/22	0/1/1/1
10	NAG	B	1	1,10	-	2/6/23/26	0/1/1/1
10	NAG	B	2	10	-	2/6/23/26	0/1/1/1
10	BMA	B	3	10	-	2/2/19/22	0/1/1/1
10	NAG	E	1	1,10	-	3/6/23/26	0/1/1/1
10	NAG	E	2	10	-	2/6/23/26	0/1/1/1
10	BMA	E	3	10	-	0/2/19/22	0/1/1/1
9	NAG	F	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	F	2	9	-	2/6/23/26	0/1/1/1
9	NAG	G	1	9,1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	G	2	9	-	2/6/23/26	0/1/1/1
11	NAG	I	1	11,1	-	0/6/23/26	0/1/1/1
11	NAG	I	2	11	-	2/6/23/26	0/1/1/1
11	BMA	I	3	11	-	0/2/19/22	0/1/1/1
11	MAN	I	4	11	-	1/2/19/22	0/1/1/1
11	MAN	I	5	11	-	0/2/19/22	0/1/1/1
9	NAG	J	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	J	2	9	-	2/6/23/26	0/1/1/1
9	NAG	K	1	9,1	-	1/6/23/26	0/1/1/1
9	NAG	K	2	9	-	0/6/23/26	0/1/1/1
10	NAG	O	1	1,10	-	2/6/23/26	0/1/1/1
10	NAG	O	2	10	-	2/6/23/26	0/1/1/1
10	BMA	O	3	10	-	2/2/19/22	0/1/1/1
9	NAG	P	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	P	2	9	-	2/6/23/26	0/1/1/1
12	NAG	S	1	12,1	-	0/6/23/26	0/1/1/1
12	NAG	S	2	12	-	4/6/23/26	0/1/1/1
12	BMA	S	3	12	-	0/2/19/22	0/1/1/1
12	MAN	S	4	12	-	0/2/19/22	0/1/1/1
12	MAN	S	5	12	-	1/2/19/22	0/1/1/1
9	NAG	T	1	9,1	-	3/6/23/26	0/1/1/1
9	NAG	T	2	9	-	2/6/23/26	0/1/1/1
13	NAG	U	1	1,13	-	2/6/23/26	0/1/1/1
13	NAG	U	2	13	-	0/6/23/26	0/1/1/1
13	BMA	U	3	13	-	0/2/19/22	0/1/1/1
13	MAN	U	4	13	-	1/2/19/22	0/1/1/1
9	NAG	V	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	V	2	9	-	2/6/23/26	0/1/1/1
10	NAG	W	1	1,10	-	2/6/23/26	0/1/1/1
10	NAG	W	2	10	-	2/6/23/26	0/1/1/1
10	BMA	W	3	10	-	2/2/19/22	0/1/1/1
10	NAG	X	1	1,10	-	1/6/23/26	0/1/1/1
10	NAG	X	2	10	-	2/6/23/26	0/1/1/1
10	BMA	X	3	10	-	2/2/19/22	0/1/1/1
9	NAG	Y	1	9,1	-	3/6/23/26	0/1/1/1
9	NAG	Y	2	9	-	2/6/23/26	0/1/1/1
9	NAG	Z	1	9,1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	Z	2	9	-	2/6/23/26	0/1/1/1
11	NAG	a	1	11,1	-	2/6/23/26	0/1/1/1
11	NAG	a	2	11	-	2/6/23/26	0/1/1/1
11	BMA	a	3	11	-	1/2/19/22	0/1/1/1
11	MAN	a	4	11	-	1/2/19/22	0/1/1/1
11	MAN	a	5	11	-	0/2/19/22	0/1/1/1
9	NAG	b	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	b	2	9	-	2/6/23/26	0/1/1/1
9	NAG	e	1	9,1	-	1/6/23/26	0/1/1/1
9	NAG	e	2	9	-	2/6/23/26	0/1/1/1
10	NAG	f	1	1,10	-	2/6/23/26	0/1/1/1
10	NAG	f	2	10	-	2/6/23/26	0/1/1/1
10	BMA	f	3	10	-	2/2/19/22	0/1/1/1
9	NAG	g	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	g	2	9	-	0/6/23/26	0/1/1/1
12	NAG	i	1	12,1	-	0/6/23/26	0/1/1/1
12	NAG	i	2	12	-	4/6/23/26	0/1/1/1
12	BMA	i	3	12	-	0/2/19/22	0/1/1/1
12	MAN	i	4	12	-	0/2/19/22	0/1/1/1
12	MAN	i	5	12	-	1/2/19/22	0/1/1/1
9	NAG	j	1	9,1	-	3/6/23/26	0/1/1/1
9	NAG	j	2	9	-	2/6/23/26	0/1/1/1
13	NAG	k	1	1,13	-	2/6/23/26	0/1/1/1
13	NAG	k	2	13	-	2/6/23/26	0/1/1/1
13	BMA	k	3	13	-	0/2/19/22	0/1/1/1
13	MAN	k	4	13	-	2/2/19/22	0/1/1/1
9	NAG	o	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	o	2	9	-	2/6/23/26	0/1/1/1
10	NAG	p	1	1,10	-	2/6/23/26	0/1/1/1
10	NAG	p	2	10	-	0/6/23/26	0/1/1/1
10	BMA	p	3	10	-	1/2/19/22	0/1/1/1
10	NAG	s	1	1,10	-	1/6/23/26	0/1/1/1
10	NAG	s	2	10	-	0/6/23/26	0/1/1/1
10	BMA	s	3	10	-	2/2/19/22	0/1/1/1
9	NAG	t	1	9,1	-	4/6/23/26	0/1/1/1
9	NAG	t	2	9	-	2/6/23/26	0/1/1/1
9	NAG	u	1	9,1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	u	2	9	-	2/6/23/26	0/1/1/1
11	NAG	v	1	11,1	-	2/6/23/26	0/1/1/1
11	NAG	v	2	11	-	2/6/23/26	0/1/1/1
11	BMA	v	3	11	-	2/2/19/22	0/1/1/1
11	MAN	v	4	11	-	1/2/19/22	0/1/1/1
11	MAN	v	5	11	-	0/2/19/22	0/1/1/1
9	NAG	w	1	9,1	-	1/6/23/26	0/1/1/1
9	NAG	w	2	9	-	2/6/23/26	0/1/1/1
9	NAG	x	1	9,1	-	1/6/23/26	0/1/1/1
9	NAG	x	2	9	-	2/6/23/26	0/1/1/1
10	NAG	y	1	1,10	-	2/6/23/26	0/1/1/1
10	NAG	y	2	10	-	2/6/23/26	0/1/1/1
10	BMA	y	3	10	-	2/2/19/22	0/1/1/1
9	NAG	z	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	z	2	9	-	0/6/23/26	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	y	3	BMA	C1-C2	2.72	1.58	1.52
10	O	3	BMA	C1-C2	2.70	1.58	1.52
11	v	4	MAN	C1-C2	2.56	1.58	1.52
12	0	3	BMA	C2-C3	2.54	1.56	1.52
12	i	3	BMA	C2-C3	2.53	1.56	1.52
13	U	4	MAN	C1-C2	2.45	1.58	1.52
9	Y	1	NAG	C1-C2	2.43	1.55	1.52
12	S	3	BMA	C2-C3	2.42	1.56	1.52
10	f	3	BMA	C1-C2	2.40	1.57	1.52
12	S	4	MAN	C1-C2	2.31	1.57	1.52
13	AA	3	BMA	C4-C3	2.30	1.58	1.52
10	B	3	BMA	C1-C2	2.29	1.57	1.52
12	0	4	MAN	C1-C2	2.26	1.57	1.52
13	AA	4	MAN	C1-C2	2.26	1.57	1.52
12	i	4	MAN	C1-C2	2.25	1.57	1.52
12	0	3	BMA	C4-C3	2.22	1.58	1.52
13	U	3	BMA	C4-C3	2.22	1.58	1.52
11	a	4	MAN	C1-C2	2.17	1.57	1.52
13	k	3	BMA	C4-C3	2.16	1.57	1.52
12	S	3	BMA	O5-C1	-2.15	1.40	1.43
9	t	1	NAG	C1-C2	2.12	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	F	1	NAG	C1-C2	2.07	1.55	1.52
12	S	3	BMA	C4-C3	2.05	1.57	1.52
13	k	4	MAN	C1-C2	2.00	1.57	1.52

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	B	3	BMA	C1-O5-C5	5.03	118.92	112.19
12	0	4	MAN	C1-O5-C5	4.80	118.62	112.19
10	W	3	BMA	C1-O5-C5	4.68	118.46	112.19
13	AA	4	MAN	C1-O5-C5	4.55	118.29	112.19
12	i	4	MAN	C1-O5-C5	4.48	118.19	112.19
10	p	3	BMA	C1-O5-C5	4.46	118.16	112.19
13	U	4	MAN	C1-O5-C5	4.21	117.83	112.19
12	S	4	MAN	C1-O5-C5	4.15	117.75	112.19
13	k	4	MAN	C1-O5-C5	3.73	117.18	112.19
9	V	1	NAG	C2-N2-C7	3.33	127.36	122.90
9	o	1	NAG	C2-N2-C7	3.29	127.31	122.90
12	i	2	NAG	C2-N2-C7	3.20	127.19	122.90
12	0	2	NAG	C2-N2-C7	3.20	127.18	122.90
12	S	2	NAG	C2-N2-C7	3.19	127.17	122.90
10	y	3	BMA	O2-C2-C3	-3.18	103.56	110.15
9	T	1	NAG	C2-N2-C7	3.18	127.16	122.90
13	U	2	NAG	C1-O5-C5	3.16	116.42	112.19
9	j	1	NAG	C2-N2-C7	3.14	127.10	122.90
9	A	1	NAG	C2-N2-C7	3.11	127.07	122.90
10	f	3	BMA	O2-C2-C3	-3.11	103.70	110.15
9	t	1	NAG	C2-N2-C7	3.09	127.05	122.90
11	v	4	MAN	C1-O5-C5	3.08	116.31	112.19
9	9	1	NAG	C2-N2-C7	3.05	126.99	122.90
10	O	3	BMA	O2-C2-C3	-3.03	103.88	110.15
9	Y	1	NAG	C2-N2-C7	3.00	126.91	122.90
9	F	1	NAG	C2-N2-C7	2.95	126.86	122.90
10	B	3	BMA	C1-C2-C3	2.81	113.73	109.64
11	a	4	MAN	C1-O5-C5	2.81	115.95	112.19
12	0	4	MAN	O2-C2-C3	-2.77	104.41	110.15
10	W	2	NAG	C1-O5-C5	2.76	115.89	112.19
10	p	2	NAG	C1-O5-C5	2.76	115.88	112.19
13	U	1	NAG	C1-O5-C5	2.73	115.85	112.19
13	k	2	NAG	C1-O5-C5	2.70	115.80	112.19
12	i	4	MAN	O2-C2-C3	-2.69	104.58	110.15
12	i	3	BMA	C1-C2-C3	2.64	113.49	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	k	1	NAG	C1-O5-C5	2.61	115.69	112.19
12	0	3	BMA	C1-C2-C3	2.60	113.43	109.64
12	S	5	MAN	C1-O5-C5	2.56	115.62	112.19
12	0	5	MAN	C1-O5-C5	2.56	115.62	112.19
13	AA	2	NAG	C1-O5-C5	2.55	115.61	112.19
13	AA	1	NAG	C1-O5-C5	2.54	115.59	112.19
11	a	5	MAN	C1-O5-C5	2.52	115.57	112.19
11	I	4	MAN	C1-O5-C5	2.52	115.56	112.19
11	I	4	MAN	O2-C2-C3	-2.49	104.99	110.15
11	a	4	MAN	O2-C2-C3	-2.49	104.99	110.15
12	S	4	MAN	O2-C2-C3	-2.48	105.02	110.15
10	O	3	BMA	C1-O5-C5	2.45	115.47	112.19
10	f	3	BMA	C1-O5-C5	2.41	115.42	112.19
11	v	4	MAN	O2-C2-C3	-2.39	105.21	110.15
10	B	2	NAG	C1-O5-C5	2.37	115.37	112.19
9	z	1	NAG	C1-O5-C5	2.36	115.34	112.19
10	B	3	BMA	O5-C1-C2	2.33	116.35	110.79
13	AA	4	MAN	O2-C2-C3	-2.28	105.42	110.15
12	S	3	BMA	C1-C2-C3	2.27	112.95	109.64
12	S	5	MAN	O2-C2-C3	-2.27	105.46	110.15
12	i	5	MAN	O2-C2-C3	-2.25	105.49	110.15
11	I	5	MAN	O2-C2-C3	-2.24	105.52	110.15
11	a	5	MAN	O2-C2-C3	-2.23	105.54	110.15
12	0	3	BMA	C2-C3-C4	2.23	114.78	110.86
13	U	2	NAG	O4-C4-C3	2.21	115.58	110.38
13	k	4	MAN	O2-C2-C3	-2.16	105.67	110.15
11	v	5	MAN	O2-C2-C3	-2.15	105.69	110.15
12	0	5	MAN	O2-C2-C3	-2.15	105.69	110.15
11	I	5	MAN	C1-O5-C5	2.15	115.07	112.19
12	i	5	MAN	C1-O5-C5	2.14	115.05	112.19
9	e	1	NAG	C1-O5-C5	2.09	114.99	112.19
9	J	2	NAG	C1-O5-C5	2.08	114.97	112.19
11	I	1	NAG	C1-O5-C5	2.07	114.96	112.19
10	B	3	BMA	C3-C4-C5	-2.04	106.53	110.23
12	i	2	NAG	C1-C2-N2	2.03	113.63	110.43
13	AA	2	NAG	O4-C4-C3	2.02	115.14	110.38
10	W	3	BMA	O2-C2-C3	-2.02	105.96	110.15
12	S	2	NAG	C1-C2-N2	2.02	113.61	110.43
13	U	4	MAN	O2-C2-C3	-2.02	105.97	110.15

There are no chirality outliers.

All (175) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	Z	2	NAG	O5-C5-C6-O6
9	A	2	NAG	O5-C5-C6-O6
9	G	1	NAG	O5-C5-C6-O6
9	u	1	NAG	O5-C5-C6-O6
12	i	2	NAG	O5-C5-C6-O6
12	0	2	NAG	O5-C5-C6-O6
13	k	1	NAG	O5-C5-C6-O6
9	V	2	NAG	O5-C5-C6-O6
9	Z	1	NAG	O5-C5-C6-O6
9	u	2	NAG	O5-C5-C6-O6
10	E	2	NAG	O5-C5-C6-O6
12	S	2	NAG	O5-C5-C6-O6
13	U	1	NAG	O5-C5-C6-O6
10	B	3	BMA	C4-C5-C6-O6
10	X	2	NAG	O5-C5-C6-O6
13	k	4	MAN	C4-C5-C6-O6
11	I	2	NAG	O5-C5-C6-O6
9	u	2	NAG	C4-C5-C6-O6
10	E	2	NAG	C4-C5-C6-O6
12	0	2	NAG	C4-C5-C6-O6
13	k	1	NAG	C4-C5-C6-O6
12	i	2	NAG	C4-C5-C6-O6
9	g	1	NAG	O5-C5-C6-O6
9	A	2	NAG	C4-C5-C6-O6
11	I	2	NAG	C4-C5-C6-O6
10	f	3	BMA	O5-C5-C6-O6
9	Z	2	NAG	C4-C5-C6-O6
13	U	1	NAG	C4-C5-C6-O6
13	AA	4	MAN	C4-C5-C6-O6
10	y	3	BMA	O5-C5-C6-O6
13	AA	4	MAN	O5-C5-C6-O6
9	Z	1	NAG	C4-C5-C6-O6
12	S	2	NAG	C4-C5-C6-O6
9	b	1	NAG	O5-C5-C6-O6
9	T	2	NAG	C4-C5-C6-O6
9	b	2	NAG	O5-C5-C6-O6
10	O	2	NAG	O5-C5-C6-O6
10	O	3	BMA	O5-C5-C6-O6
13	AA	1	NAG	O5-C5-C6-O6
9	G	1	NAG	C4-C5-C6-O6
9	u	1	NAG	C4-C5-C6-O6
9	x	2	NAG	C4-C5-C6-O6
10	B	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
9	J	2	NAG	O5-C5-C6-O6
10	E	1	NAG	O5-C5-C6-O6
9	w	2	NAG	O5-C5-C6-O6
9	9	1	NAG	O5-C5-C6-O6
9	V	2	NAG	C4-C5-C6-O6
13	k	4	MAN	O5-C5-C6-O6
10	O	2	NAG	C4-C5-C6-O6
13	AA	1	NAG	C4-C5-C6-O6
10	B	2	NAG	O5-C5-C6-O6
10	B	3	BMA	O5-C5-C6-O6
9	g	1	NAG	C4-C5-C6-O6
10	X	2	NAG	C4-C5-C6-O6
9	b	1	NAG	C4-C5-C6-O6
9	b	2	NAG	C4-C5-C6-O6
9	w	2	NAG	C4-C5-C6-O6
10	E	1	NAG	C4-C5-C6-O6
9	9	1	NAG	C4-C5-C6-O6
9	J	2	NAG	C4-C5-C6-O6
10	B	1	NAG	O5-C5-C6-O6
9	J	1	NAG	O5-C5-C6-O6
9	o	2	NAG	O5-C5-C6-O6
10	W	1	NAG	O5-C5-C6-O6
11	v	2	NAG	O5-C5-C6-O6
10	p	1	NAG	O5-C5-C6-O6
12	i	5	MAN	O5-C5-C6-O6
10	s	3	BMA	O5-C5-C6-O6
12	S	5	MAN	O5-C5-C6-O6
13	U	4	MAN	O5-C5-C6-O6
9	o	2	NAG	C4-C5-C6-O6
11	v	2	NAG	C4-C5-C6-O6
10	s	3	BMA	C4-C5-C6-O6
9	T	2	NAG	O5-C5-C6-O6
9	x	2	NAG	O5-C5-C6-O6
13	AA	2	NAG	C4-C5-C6-O6
13	AA	2	NAG	O5-C5-C6-O6
9	F	2	NAG	O5-C5-C6-O6
9	z	1	NAG	C4-C5-C6-O6
10	W	3	BMA	C4-C5-C6-O6
9	F	2	NAG	C4-C5-C6-O6
9	G	2	NAG	O5-C5-C6-O6
12	0	5	MAN	O5-C5-C6-O6
9	G	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
11	v	3	BMA	C4-C5-C6-O6
9	P	2	NAG	C4-C5-C6-O6
10	f	3	BMA	C4-C5-C6-O6
13	k	2	NAG	C4-C5-C6-O6
11	a	4	MAN	O5-C5-C6-O6
13	k	2	NAG	O5-C5-C6-O6
10	s	1	NAG	O5-C5-C6-O6
11	I	4	MAN	O5-C5-C6-O6
11	v	4	MAN	O5-C5-C6-O6
9	T	1	NAG	O5-C5-C6-O6
10	X	1	NAG	O5-C5-C6-O6
10	f	1	NAG	C4-C5-C6-O6
9	j	1	NAG	O5-C5-C6-O6
9	w	1	NAG	O5-C5-C6-O6
9	x	1	NAG	O5-C5-C6-O6
9	K	1	NAG	O5-C5-C6-O6
10	O	3	BMA	C4-C5-C6-O6
9	e	1	NAG	O5-C5-C6-O6
9	z	1	NAG	O5-C5-C6-O6
10	y	3	BMA	C4-C5-C6-O6
9	P	2	NAG	O5-C5-C6-O6
10	O	1	NAG	C4-C5-C6-O6
9	Y	2	NAG	C4-C5-C6-O6
9	j	2	NAG	C4-C5-C6-O6
10	W	3	BMA	O5-C5-C6-O6
9	A	1	NAG	C1-C2-N2-C7
9	V	1	NAG	C1-C2-N2-C7
9	j	1	NAG	C1-C2-N2-C7
9	o	1	NAG	C1-C2-N2-C7
9	A	1	NAG	C4-C5-C6-O6
9	e	2	NAG	C4-C5-C6-O6
10	W	2	NAG	C4-C5-C6-O6
9	Y	2	NAG	O5-C5-C6-O6
11	v	3	BMA	O5-C5-C6-O6
10	B	1	NAG	C4-C5-C6-O6
9	J	1	NAG	C4-C5-C6-O6
10	f	1	NAG	O5-C5-C6-O6
11	a	1	NAG	C4-C5-C6-O6
9	o	1	NAG	C4-C5-C6-O6
10	y	1	NAG	C4-C5-C6-O6
10	W	1	NAG	C4-C5-C6-O6
9	t	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
11	v	1	NAG	C4-C5-C6-O6
9	F	1	NAG	C4-C5-C6-O6
10	y	2	NAG	C4-C5-C6-O6
9	A	1	NAG	O5-C5-C6-O6
10	O	1	NAG	O5-C5-C6-O6
9	t	1	NAG	C4-C5-C6-O6
10	p	1	NAG	C4-C5-C6-O6
11	a	3	BMA	C4-C5-C6-O6
10	X	3	BMA	C4-C5-C6-O6
9	t	2	NAG	O5-C5-C6-O6
9	e	2	NAG	O5-C5-C6-O6
11	a	2	NAG	C4-C5-C6-O6
9	Y	1	NAG	C4-C5-C6-O6
10	y	2	NAG	O5-C5-C6-O6
10	W	2	NAG	O5-C5-C6-O6
10	X	3	BMA	O5-C5-C6-O6
10	f	2	NAG	C4-C5-C6-O6
9	o	1	NAG	O5-C5-C6-O6
10	y	1	NAG	O5-C5-C6-O6
9	F	1	NAG	C1-C2-N2-C7
9	T	1	NAG	C1-C2-N2-C7
9	Y	1	NAG	C1-C2-N2-C7
9	t	1	NAG	C1-C2-N2-C7
9	9	1	NAG	C1-C2-N2-C7
10	E	1	NAG	C1-C2-N2-C7
12	S	2	NAG	C1-C2-N2-C7
12	i	2	NAG	C1-C2-N2-C7
12	0	2	NAG	C1-C2-N2-C7
9	j	2	NAG	O5-C5-C6-O6
9	A	1	NAG	C3-C2-N2-C7
9	F	1	NAG	C3-C2-N2-C7
9	T	1	NAG	C3-C2-N2-C7
9	V	1	NAG	C3-C2-N2-C7
9	Y	1	NAG	C3-C2-N2-C7
9	j	1	NAG	C3-C2-N2-C7
9	o	1	NAG	C3-C2-N2-C7
9	t	1	NAG	C3-C2-N2-C7
9	9	1	NAG	C3-C2-N2-C7
12	S	2	NAG	C3-C2-N2-C7
12	i	2	NAG	C3-C2-N2-C7
12	0	2	NAG	C3-C2-N2-C7
10	p	3	BMA	C4-C5-C6-O6

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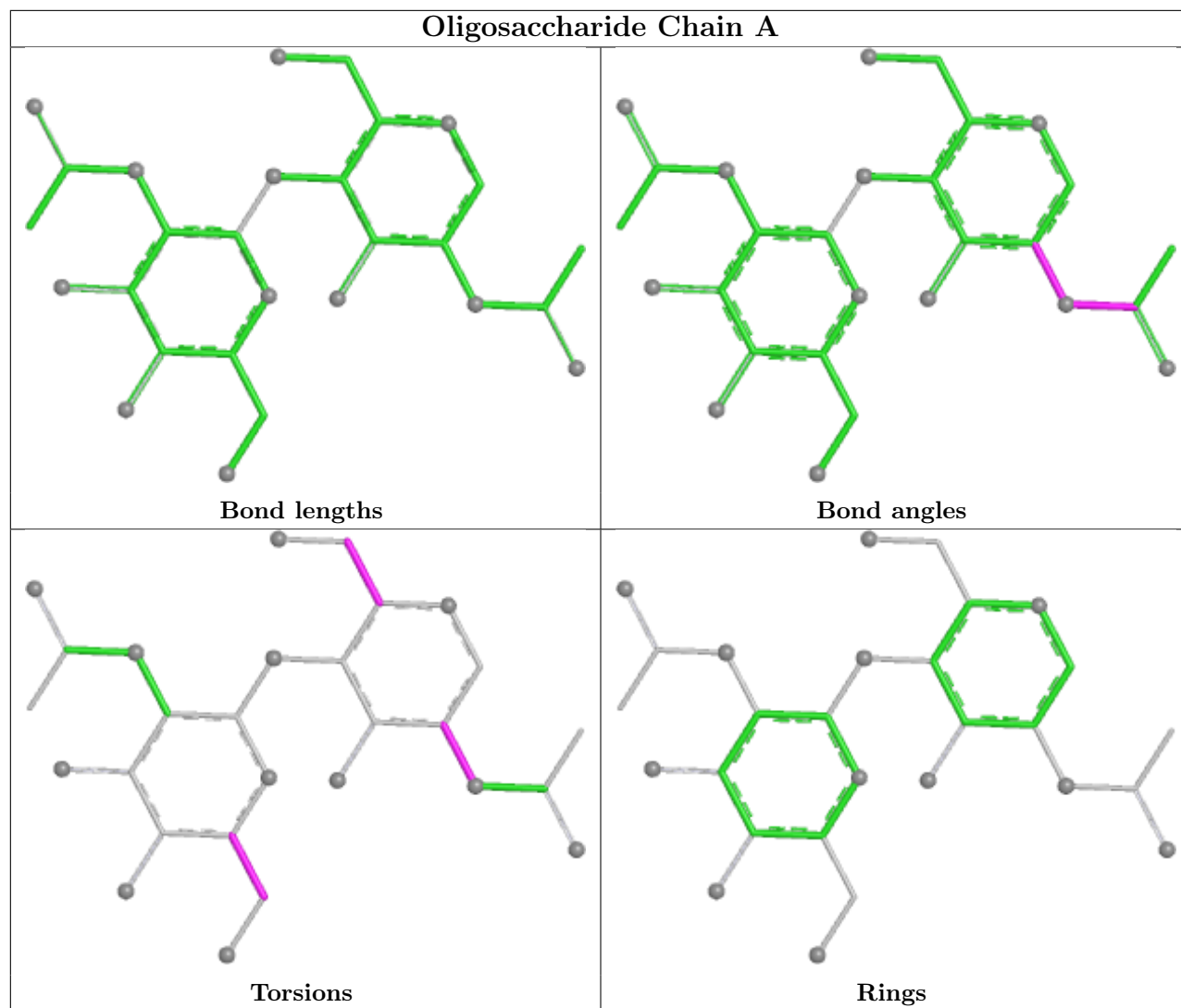
Mol	Chain	Res	Type	Atoms
11	a	2	NAG	O5-C5-C6-O6
10	f	2	NAG	O5-C5-C6-O6
11	a	1	NAG	O5-C5-C6-O6
9	F	1	NAG	O5-C5-C6-O6
9	t	1	NAG	O5-C5-C6-O6
11	v	1	NAG	O5-C5-C6-O6

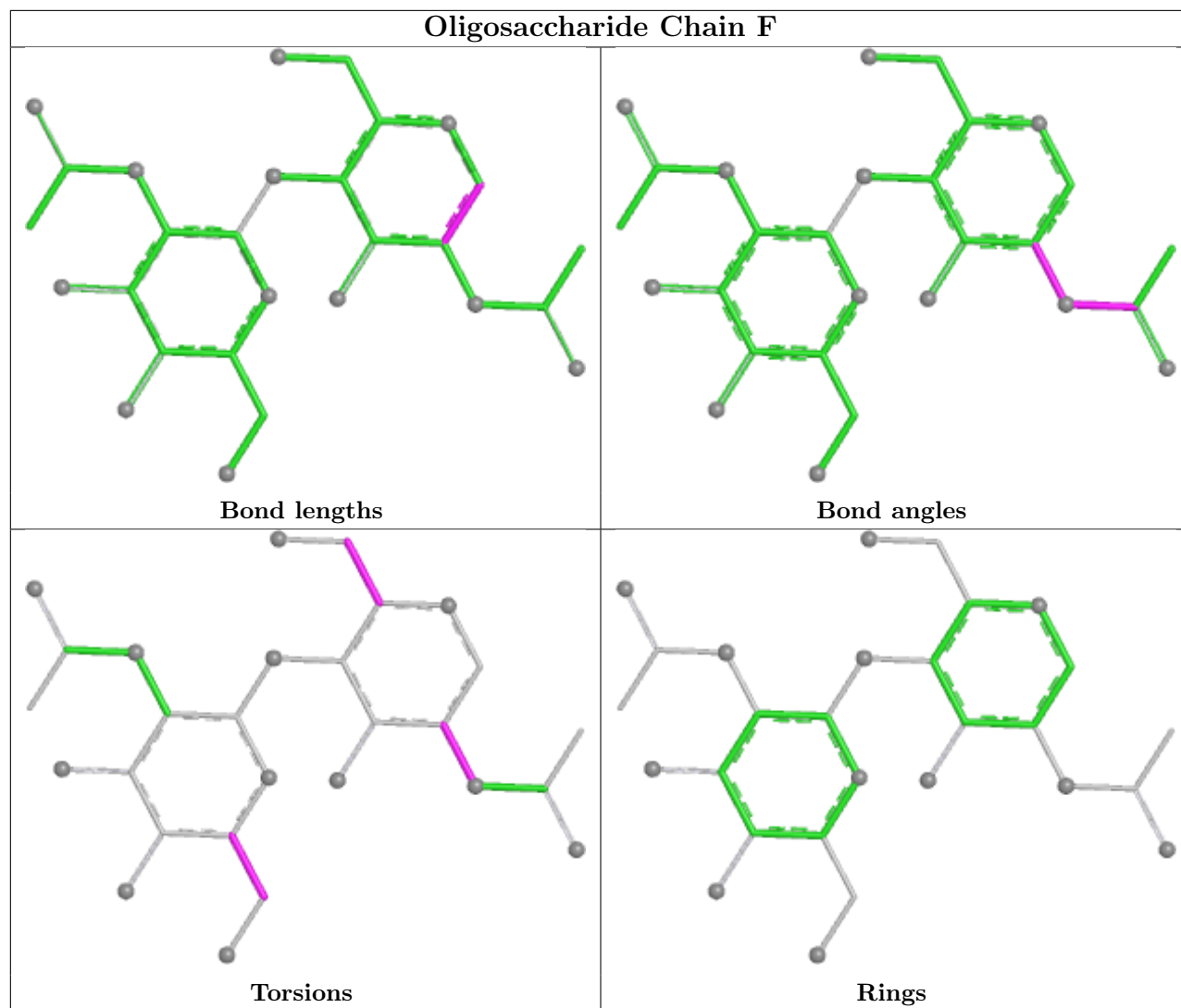
There are no ring outliers.

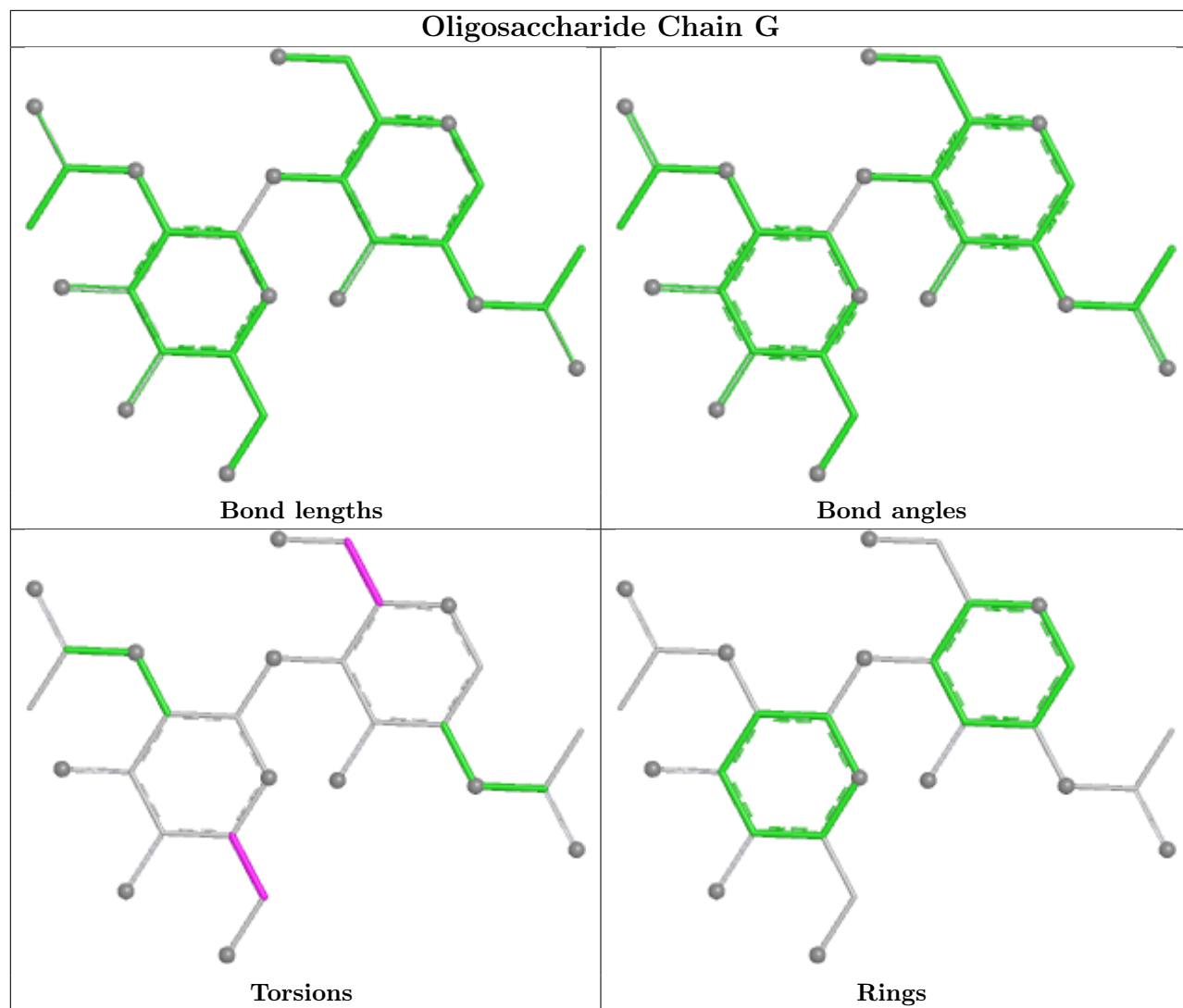
8 monomers are involved in 8 short contacts:

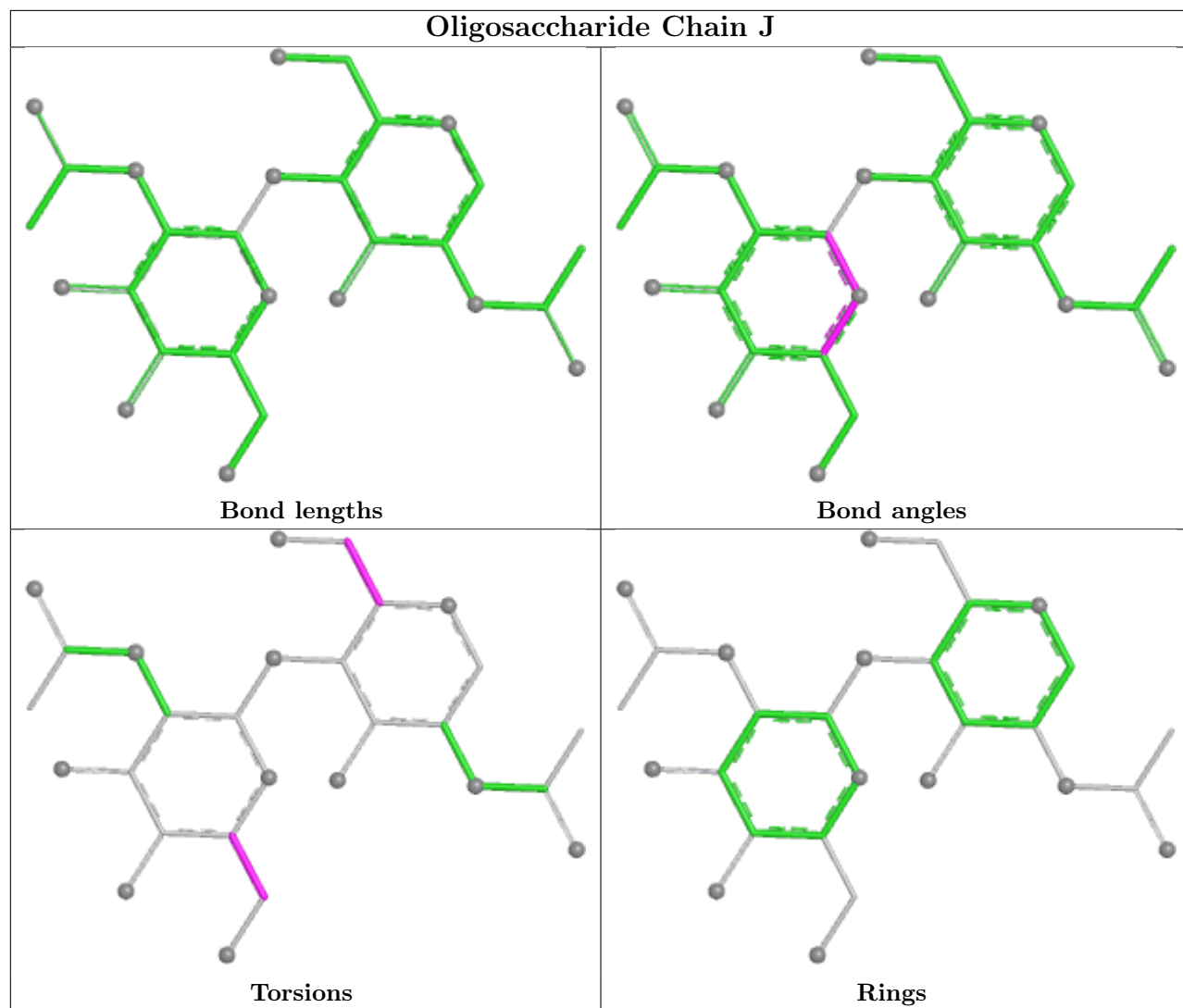
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	t	1	NAG	1	0
10	W	1	NAG	1	0
12	S	4	MAN	1	0
10	B	1	NAG	1	0
9	Y	1	NAG	1	0
13	k	1	NAG	1	0
12	i	4	MAN	1	0
10	p	1	NAG	1	0

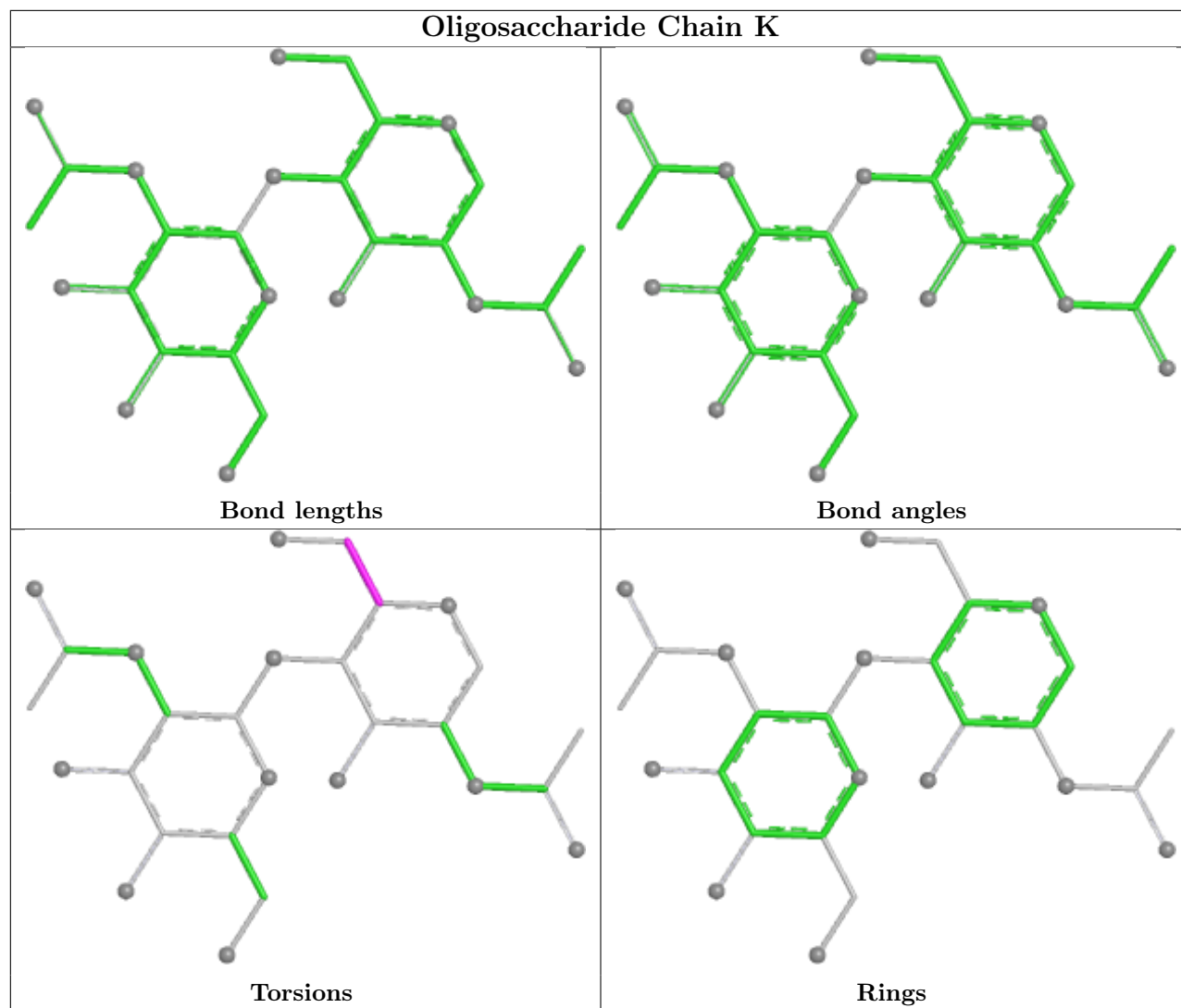
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

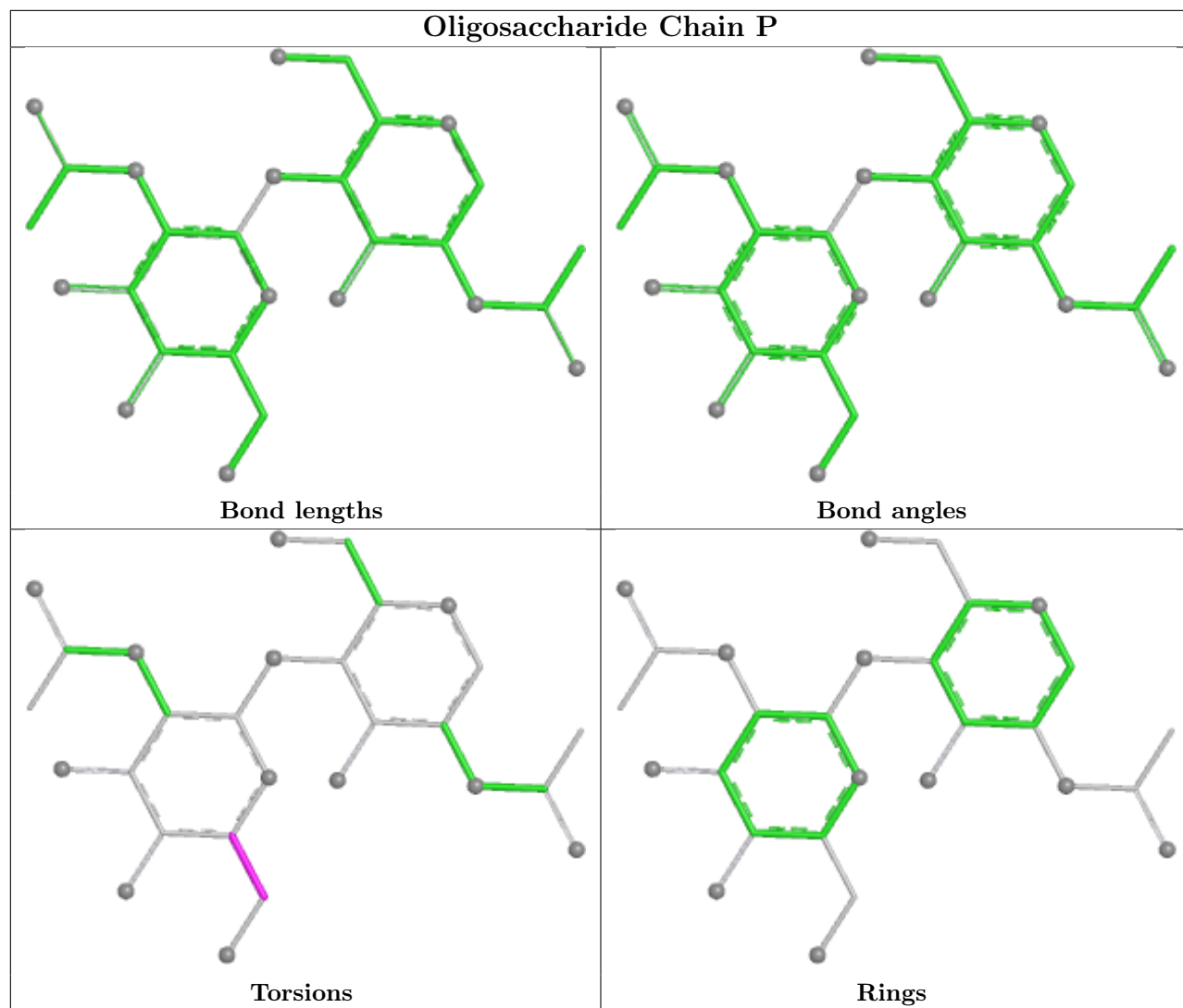


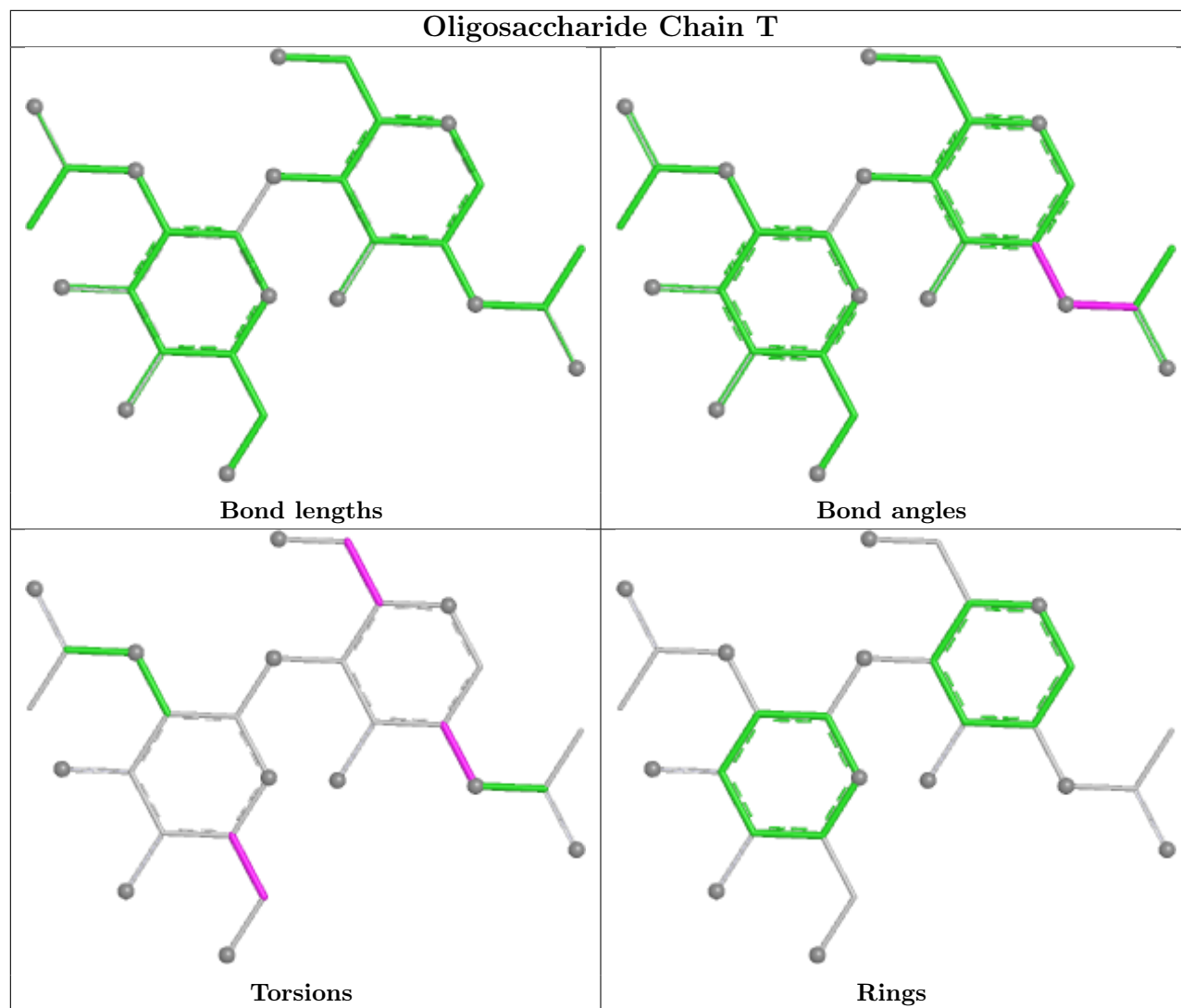


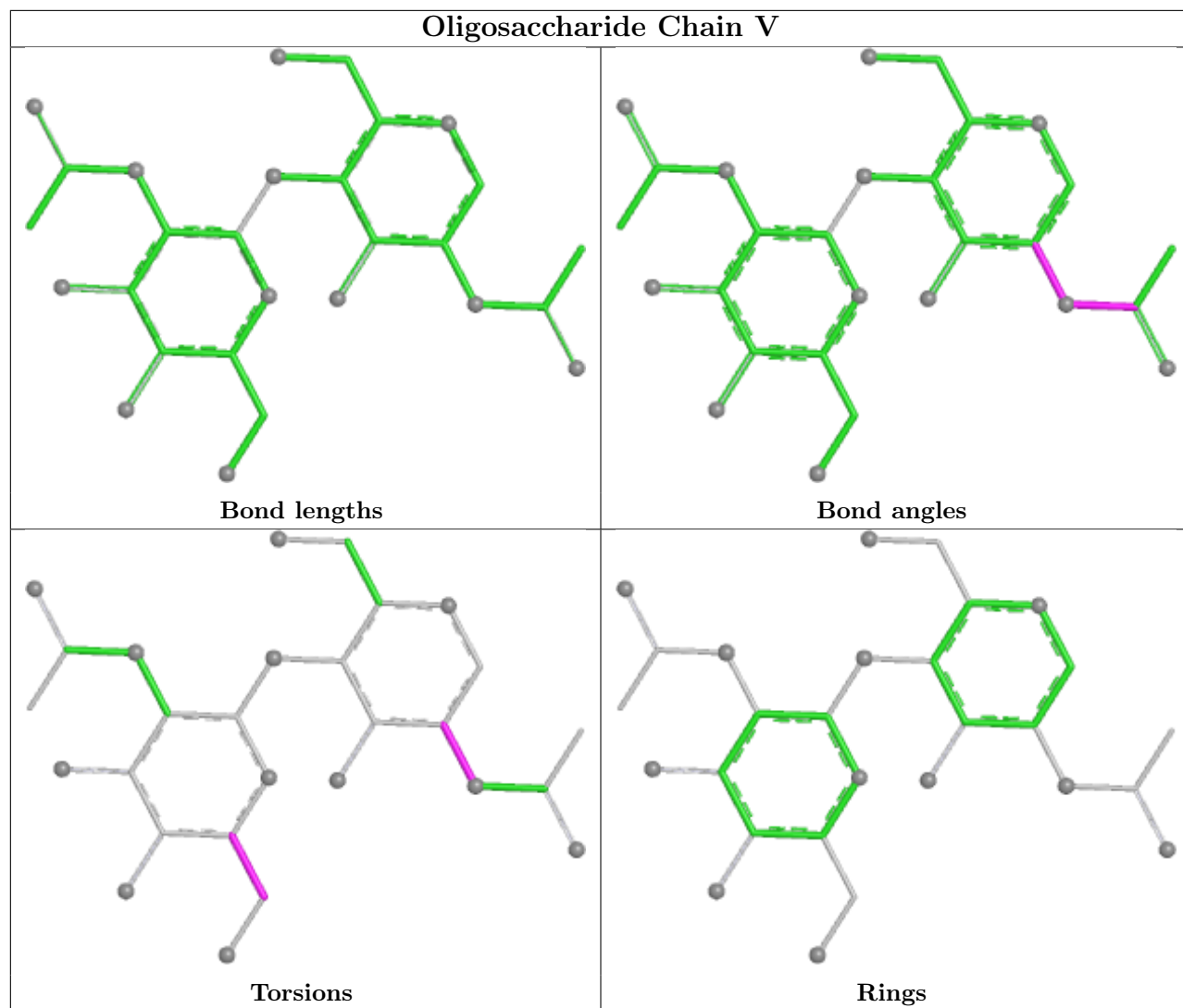


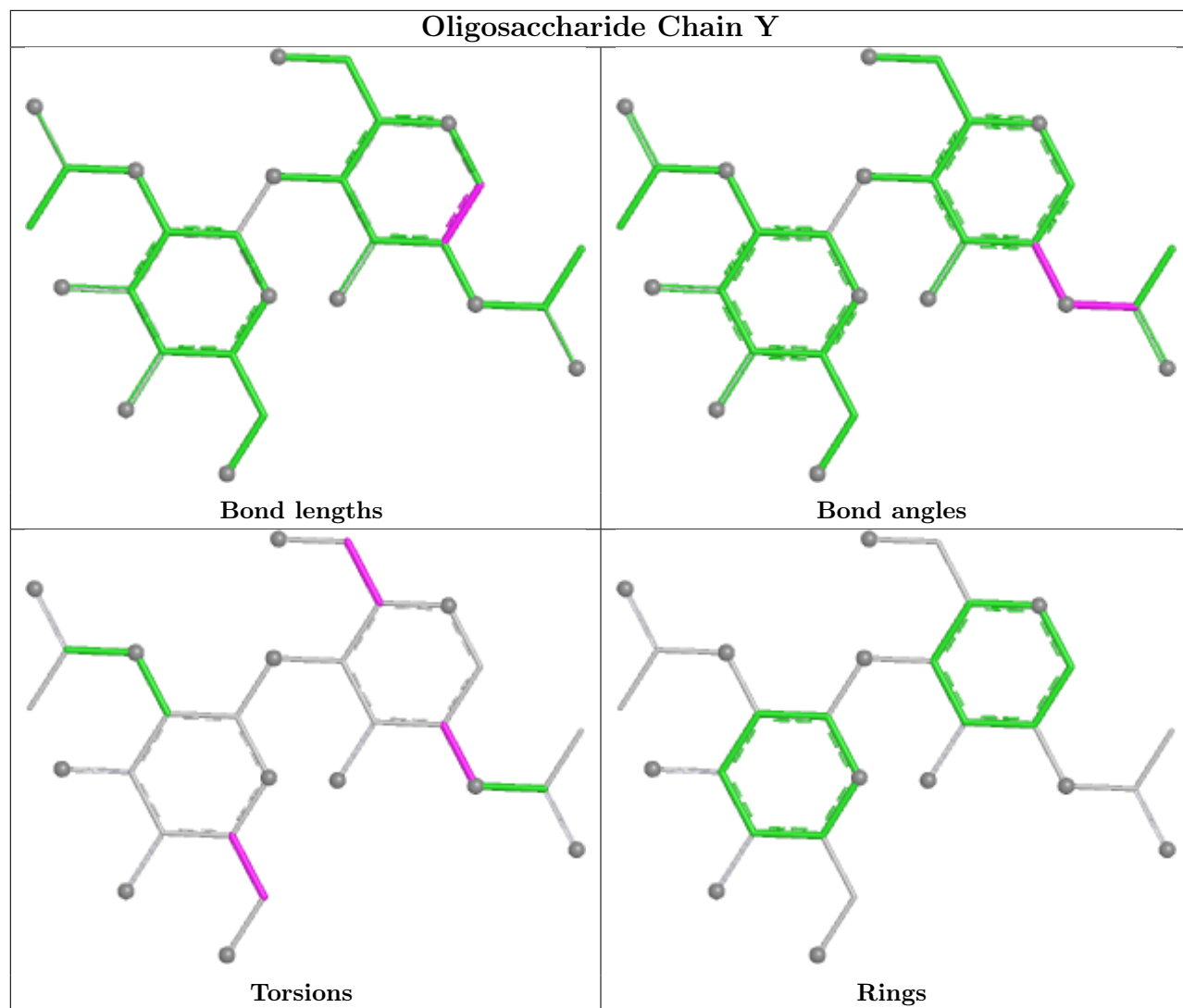


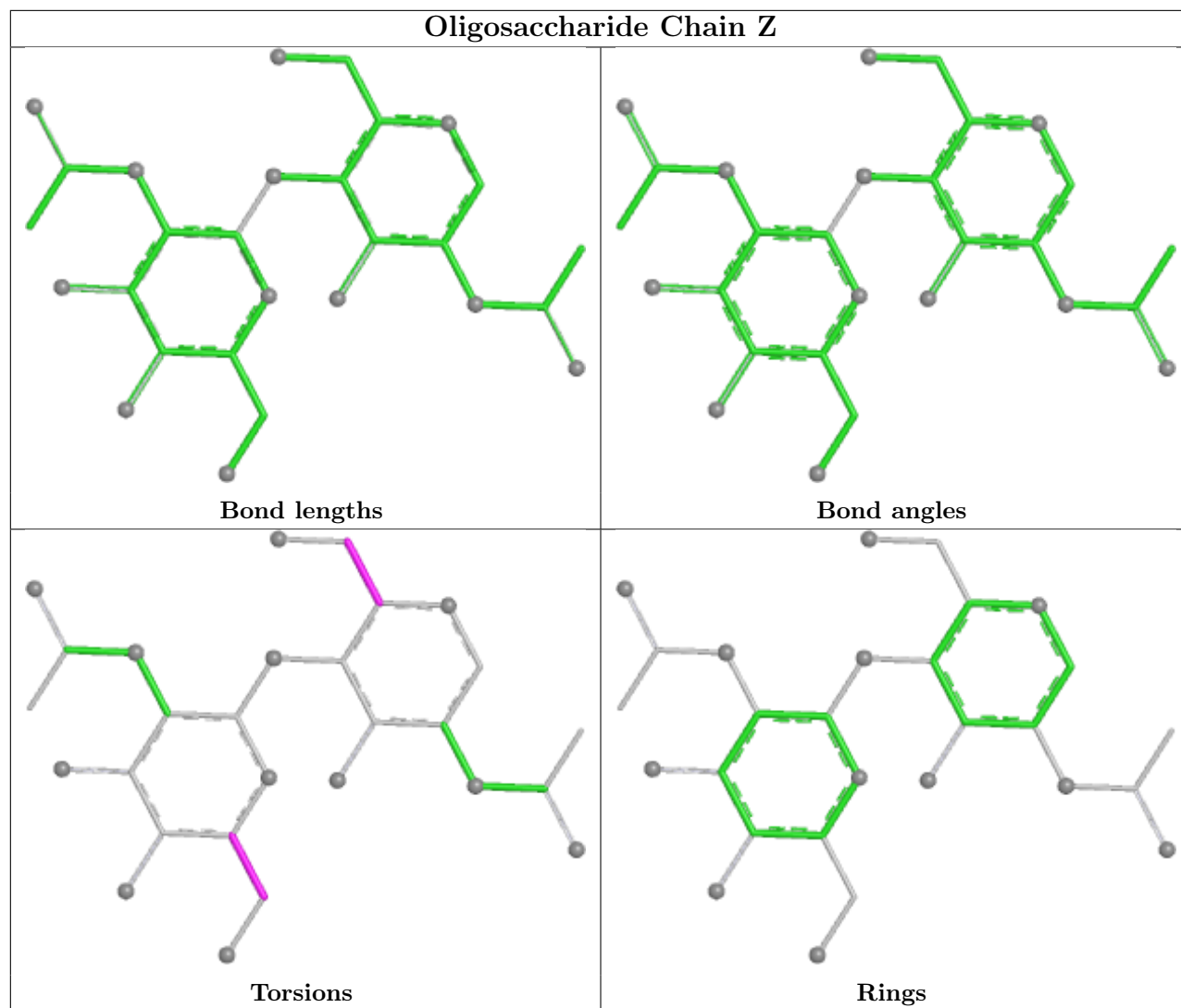


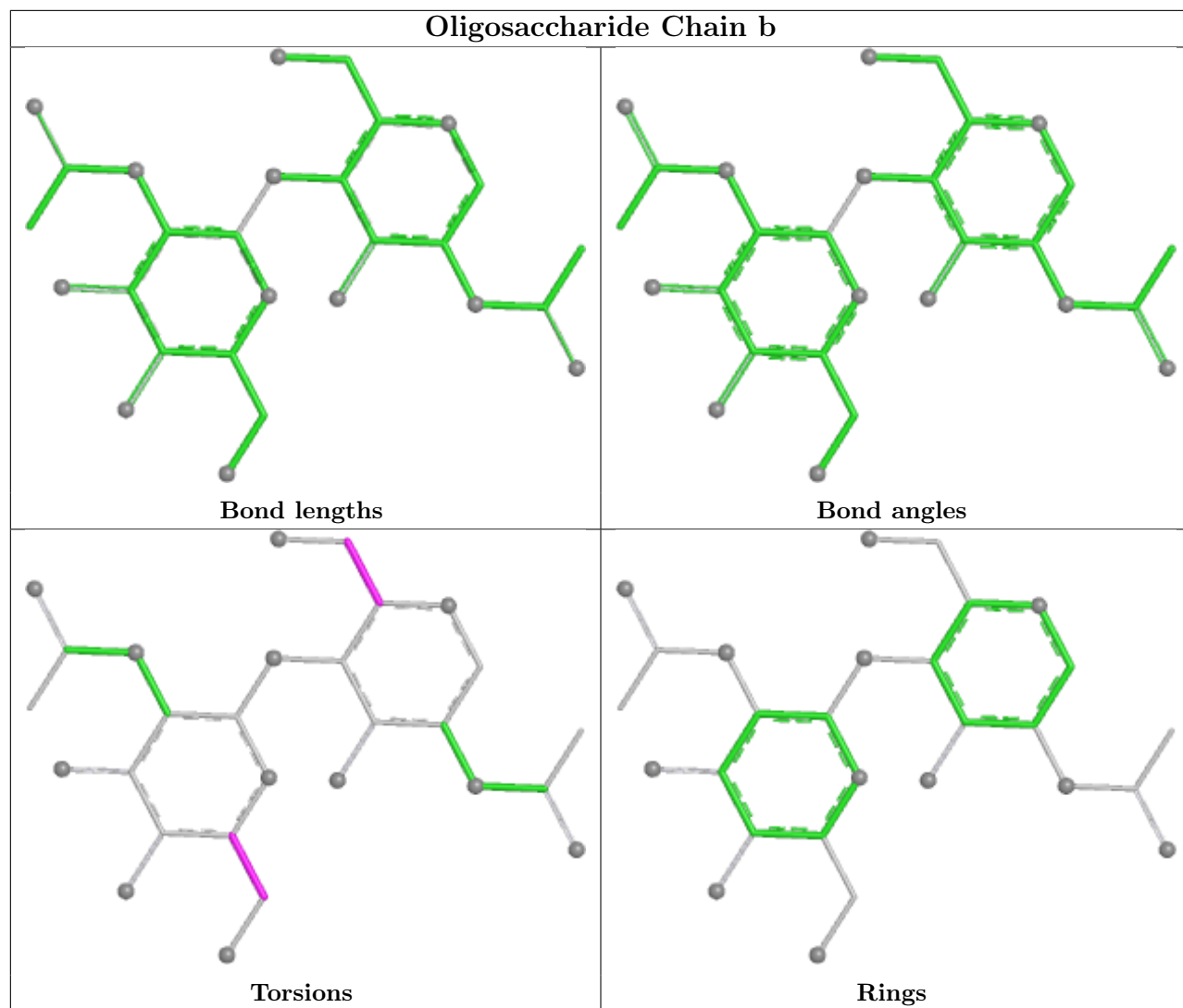


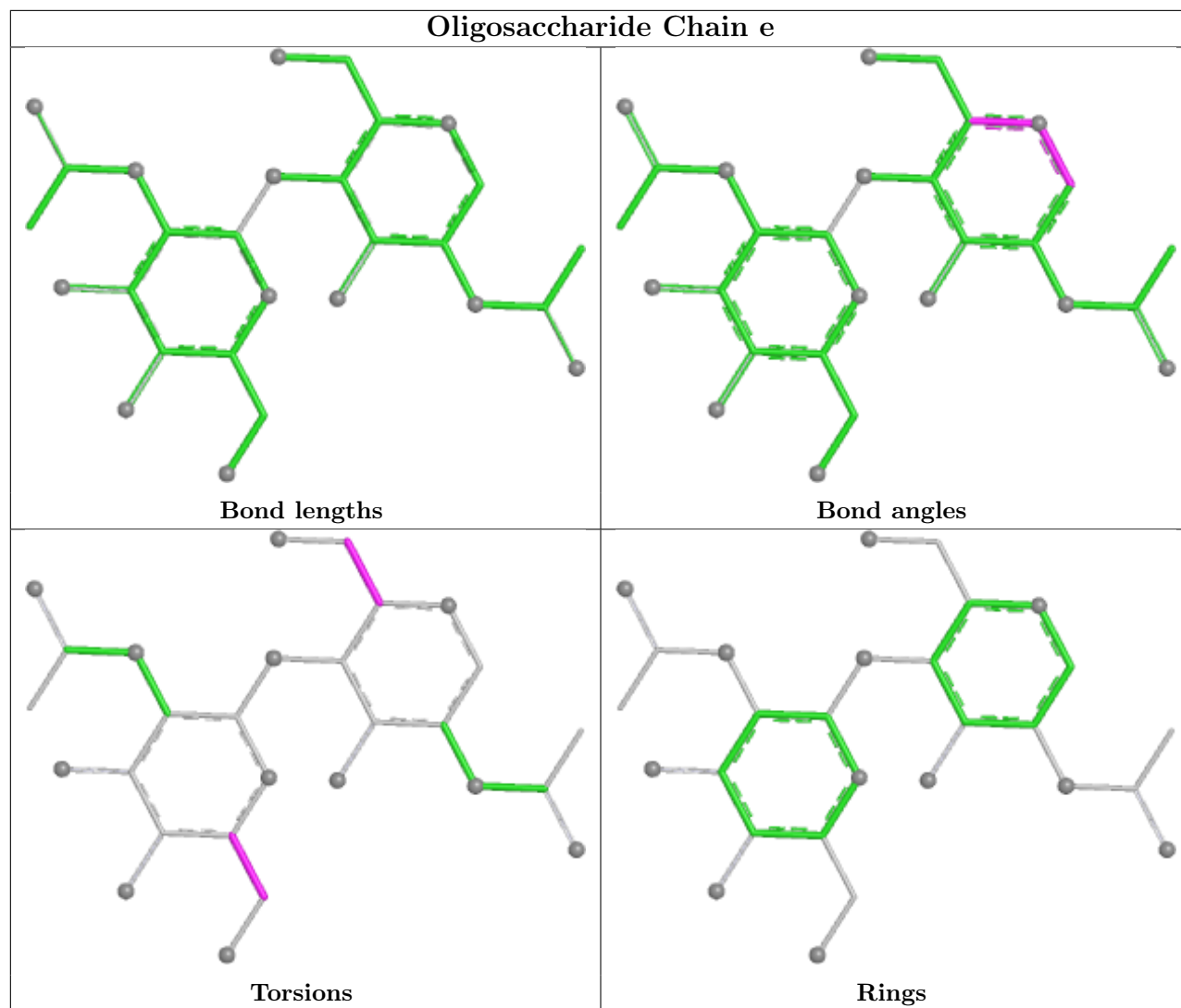


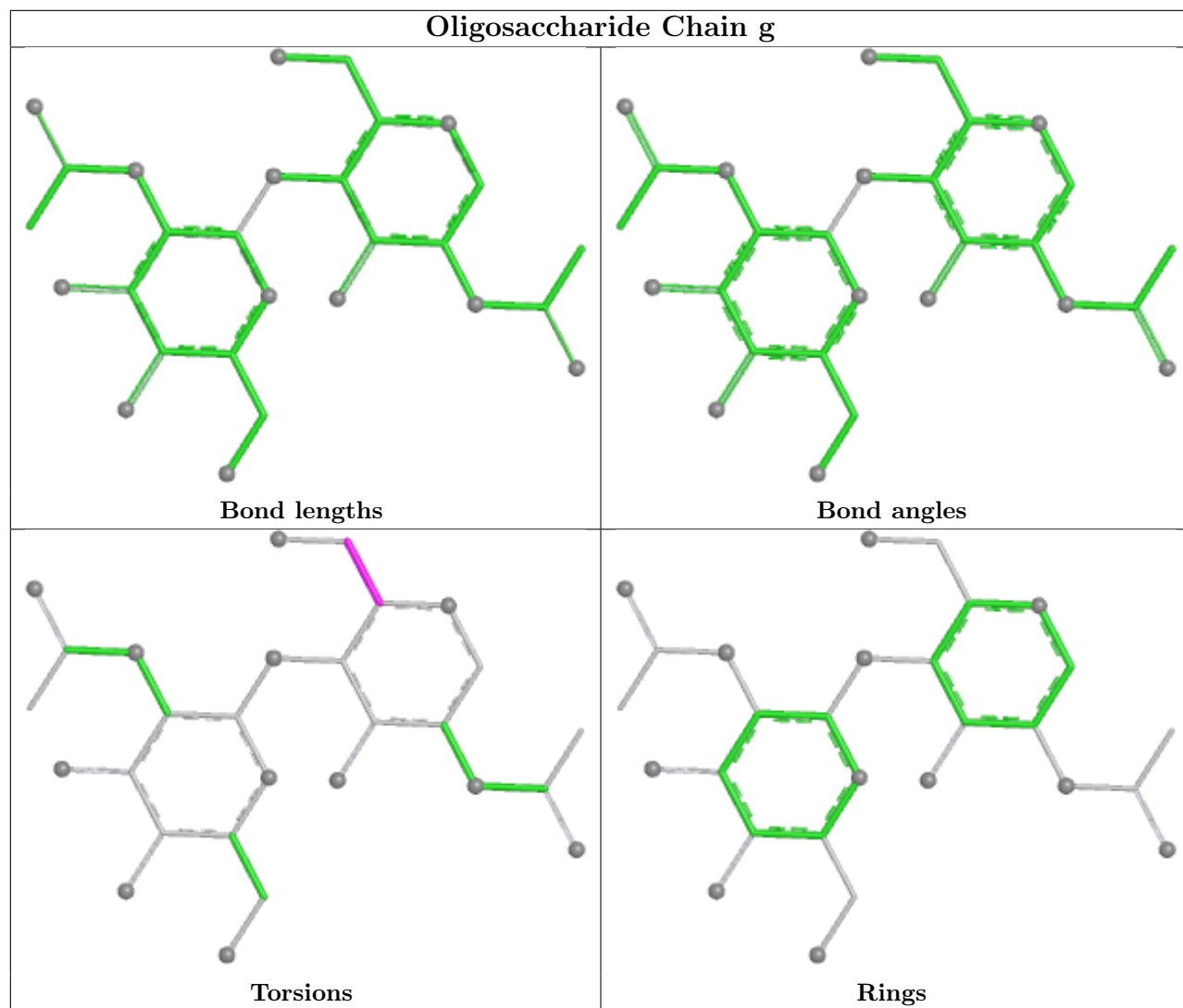


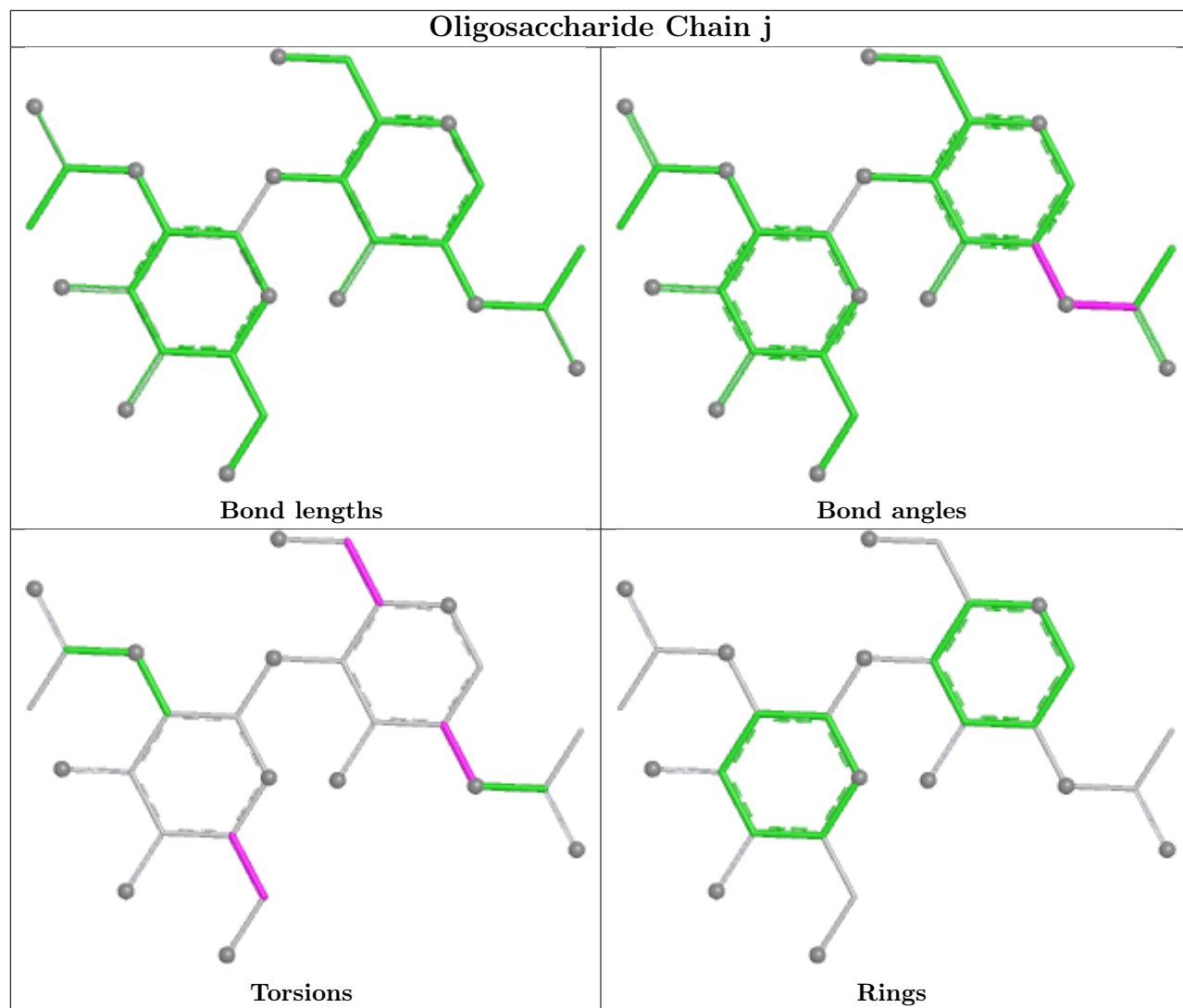


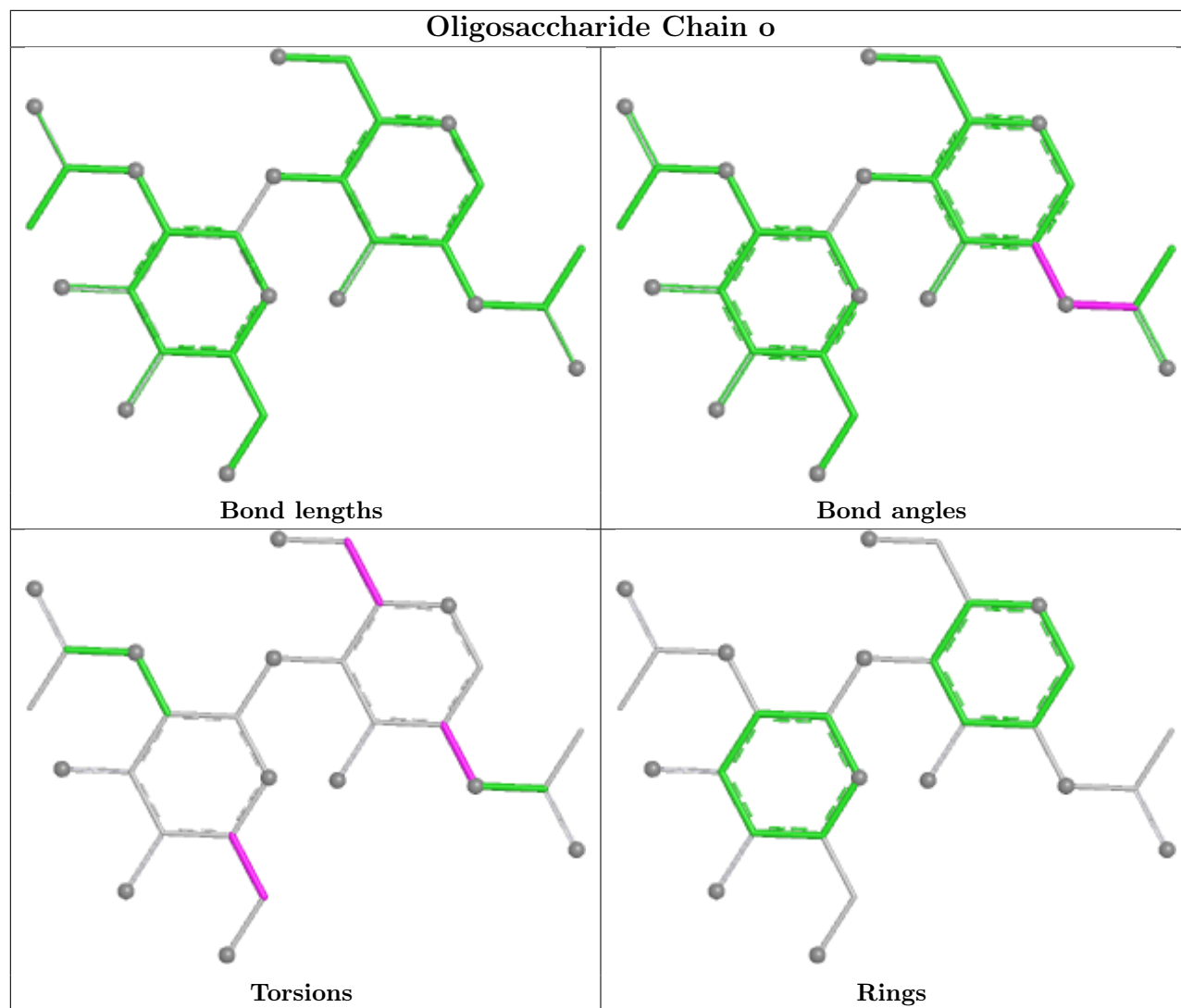


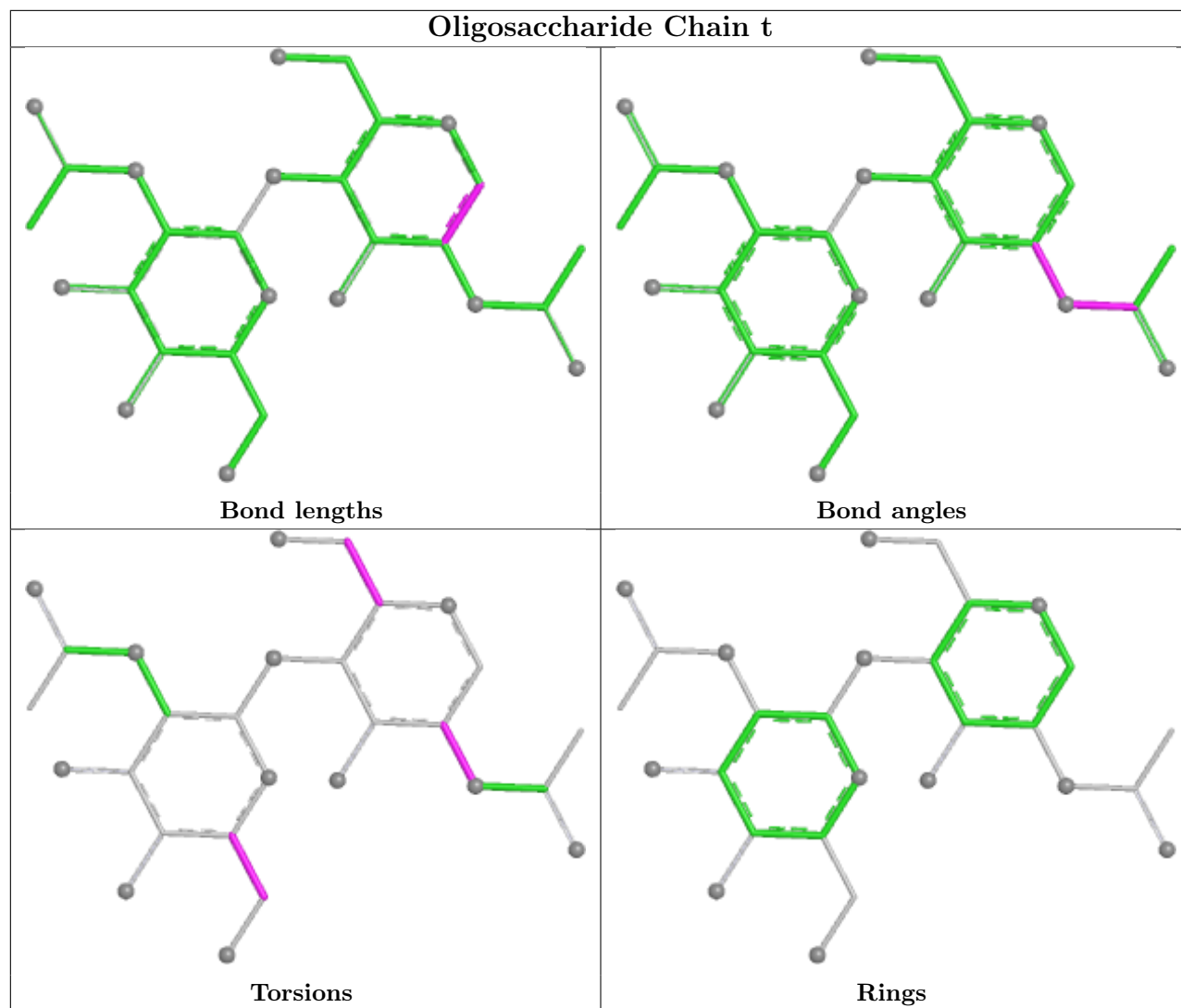


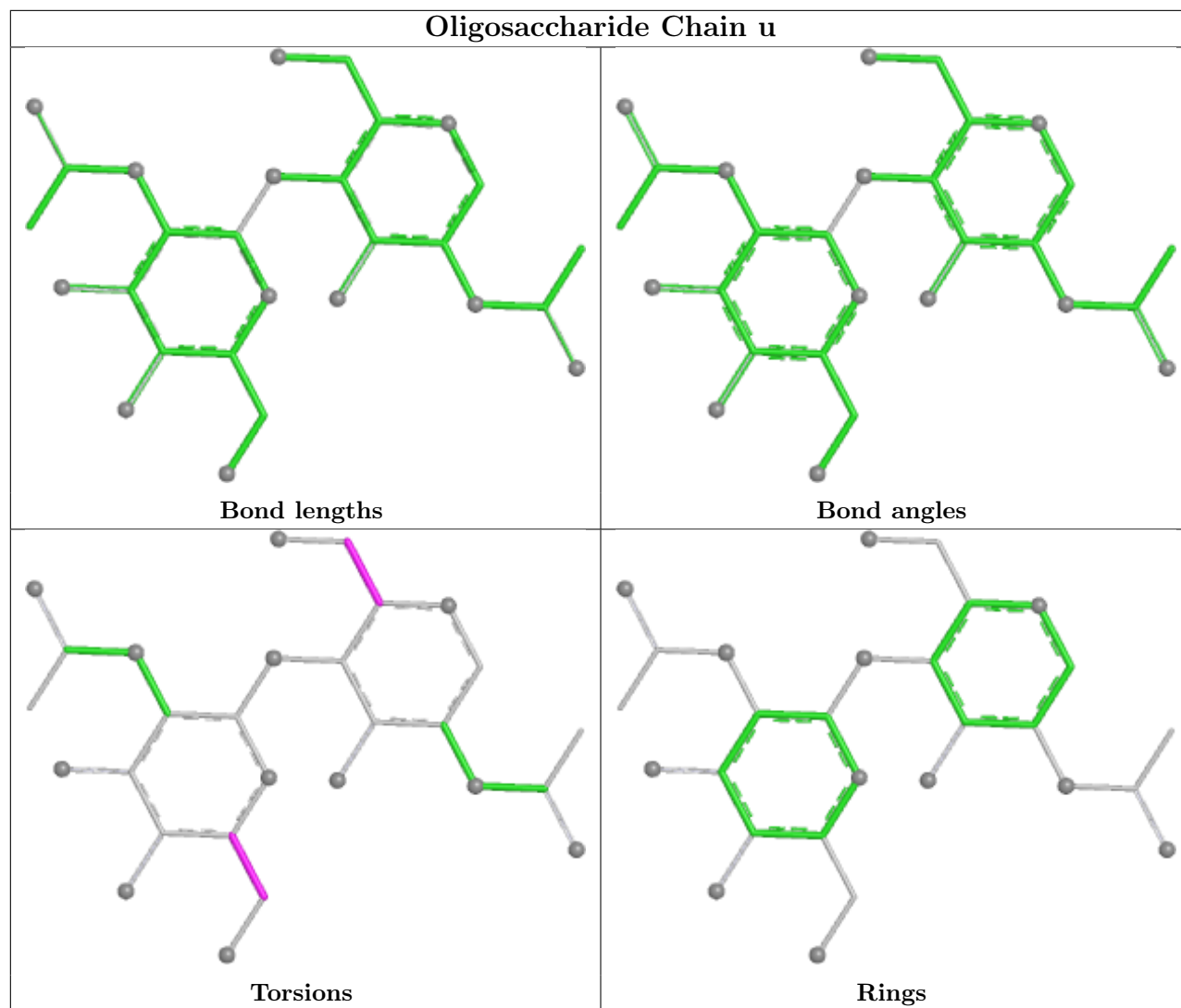


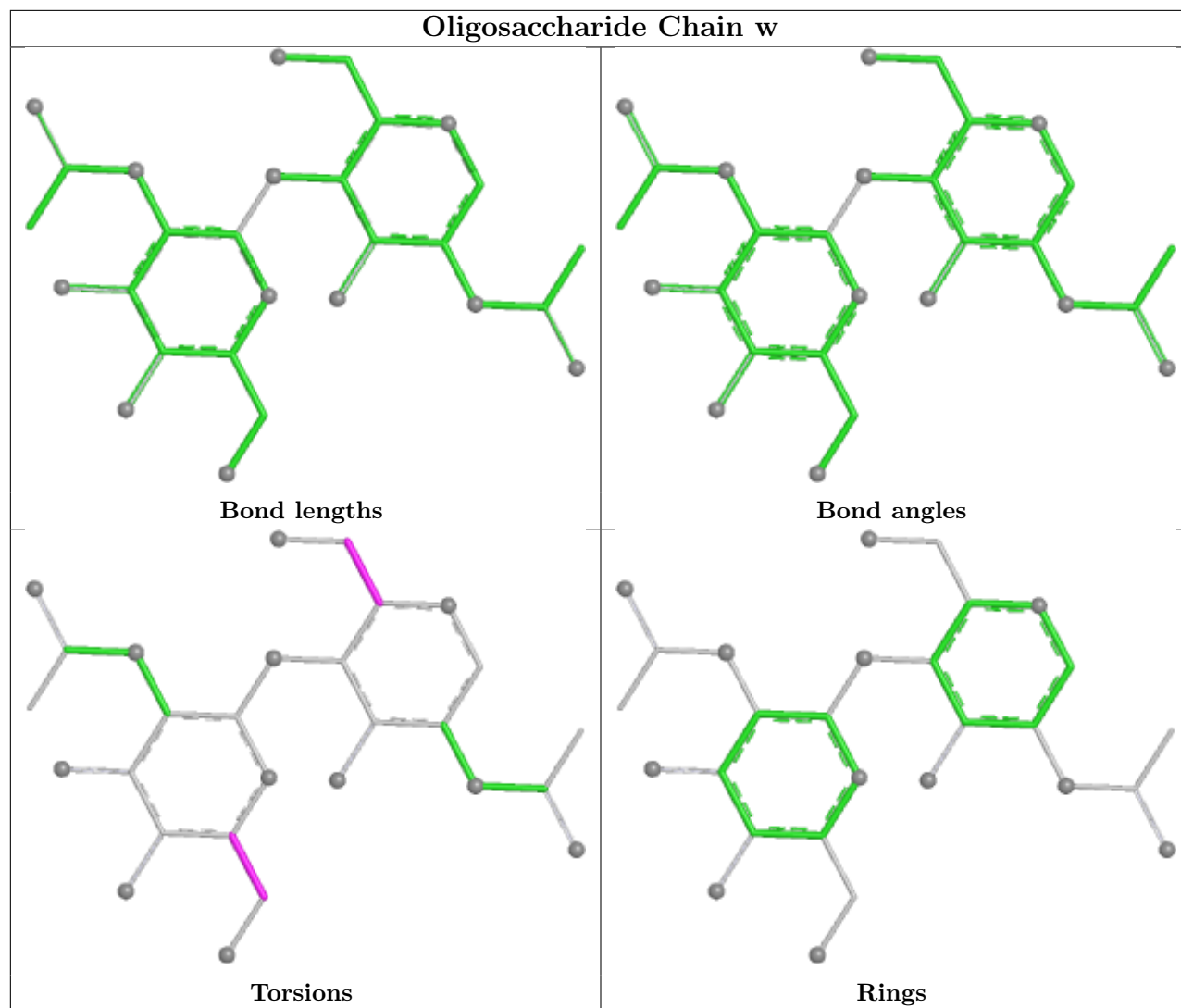


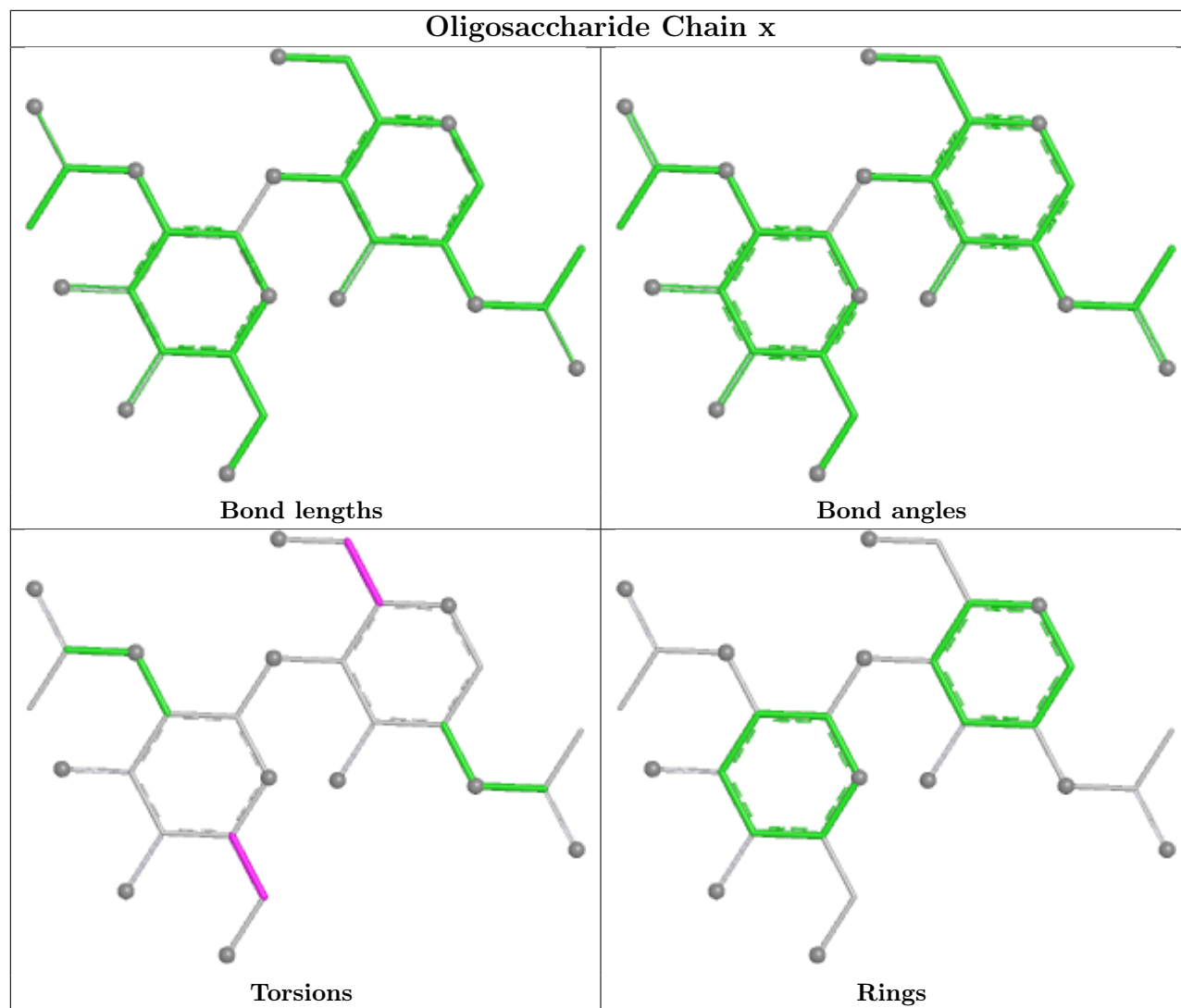


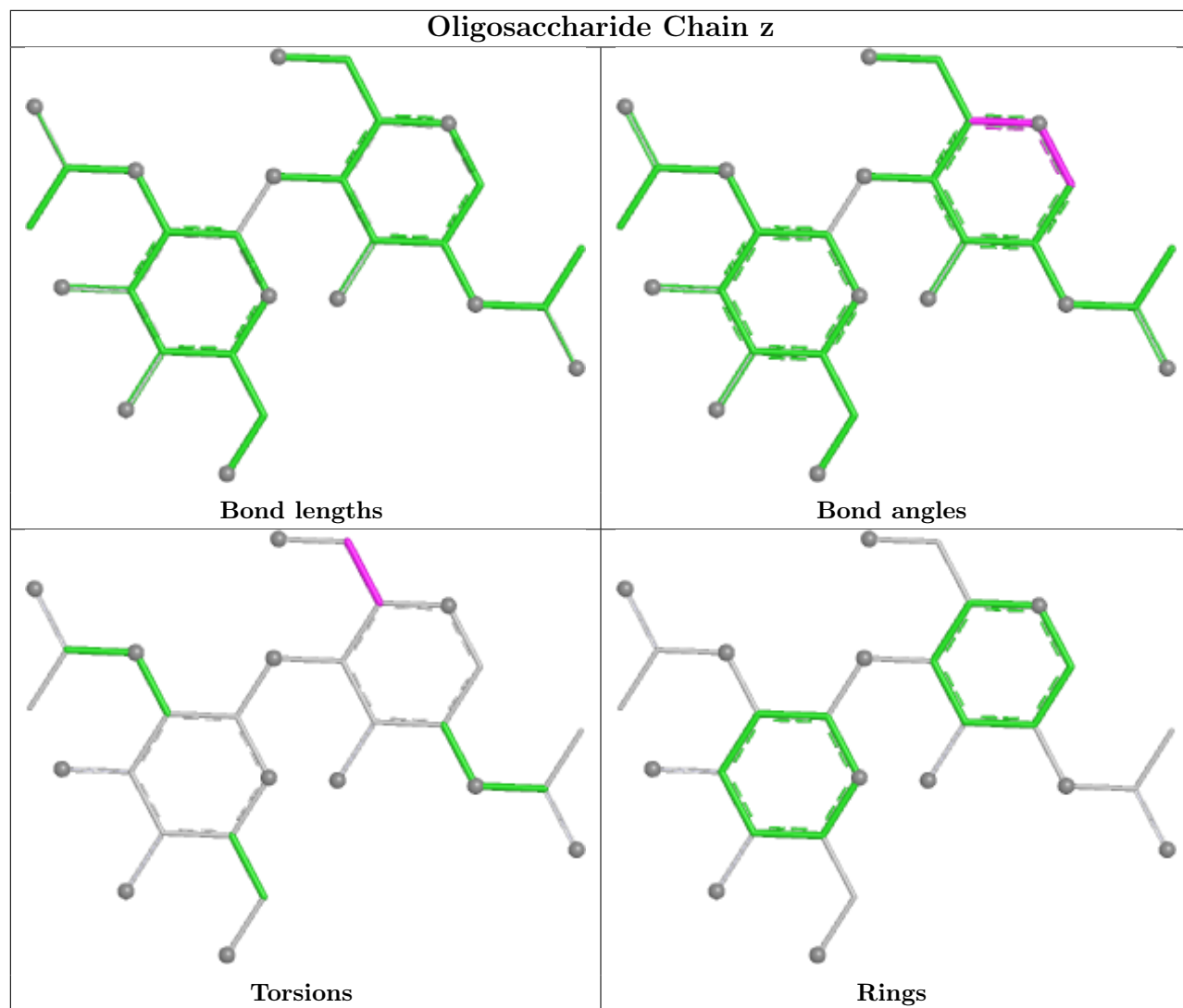


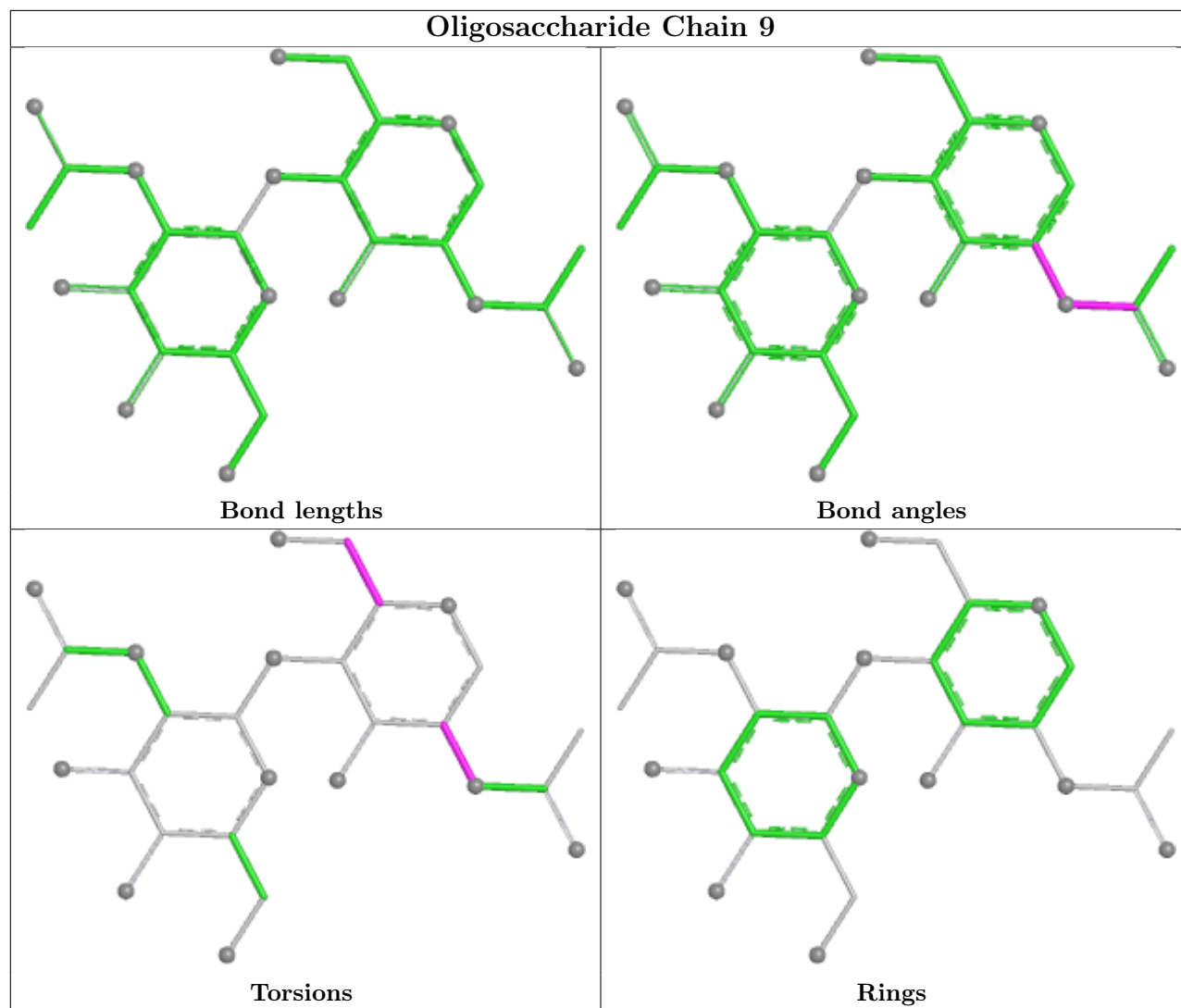


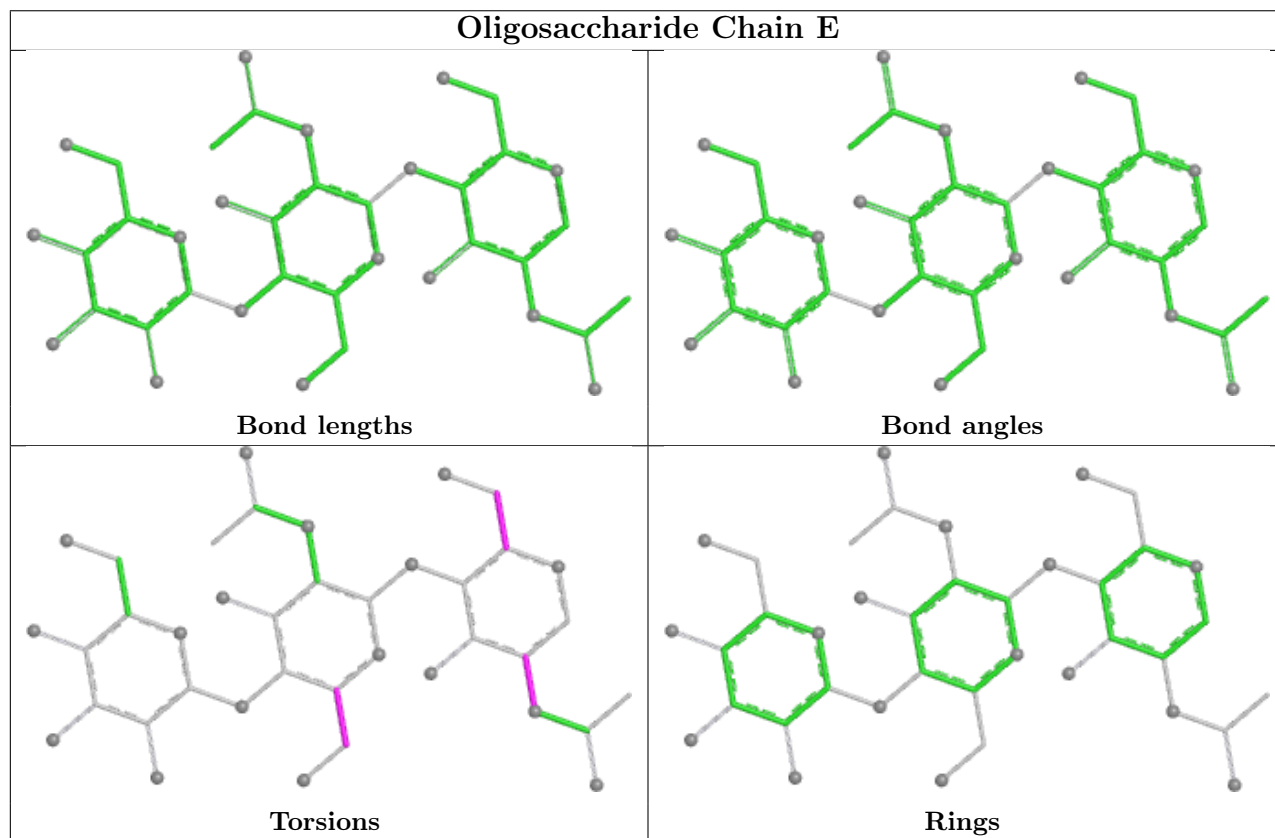
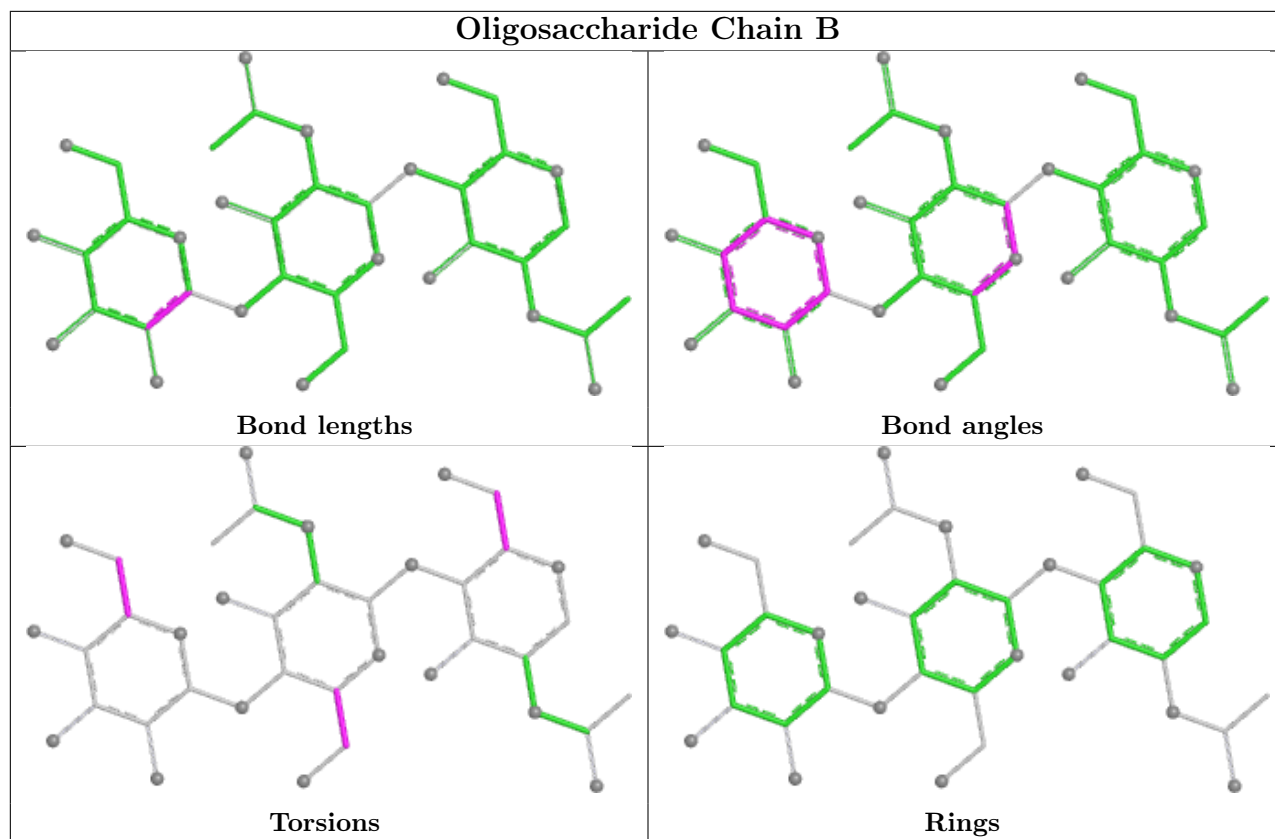


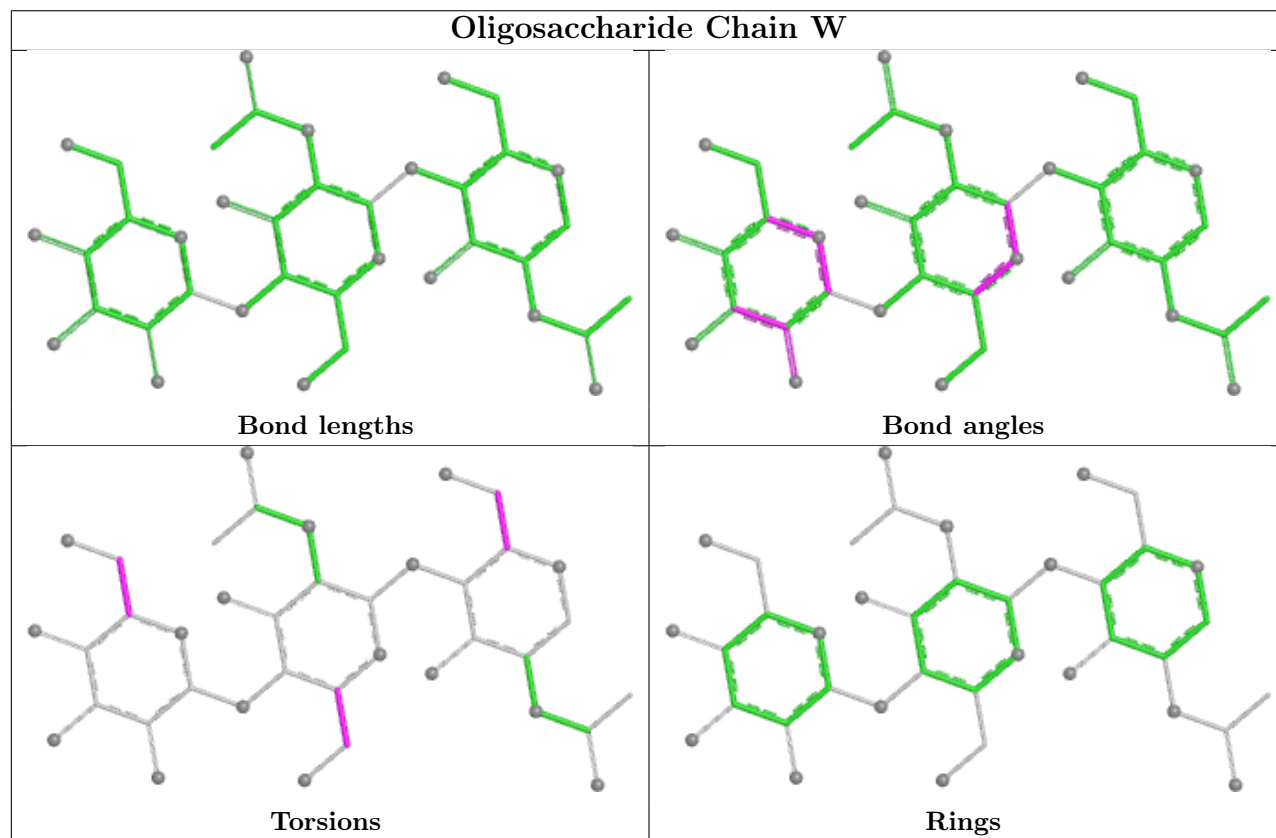
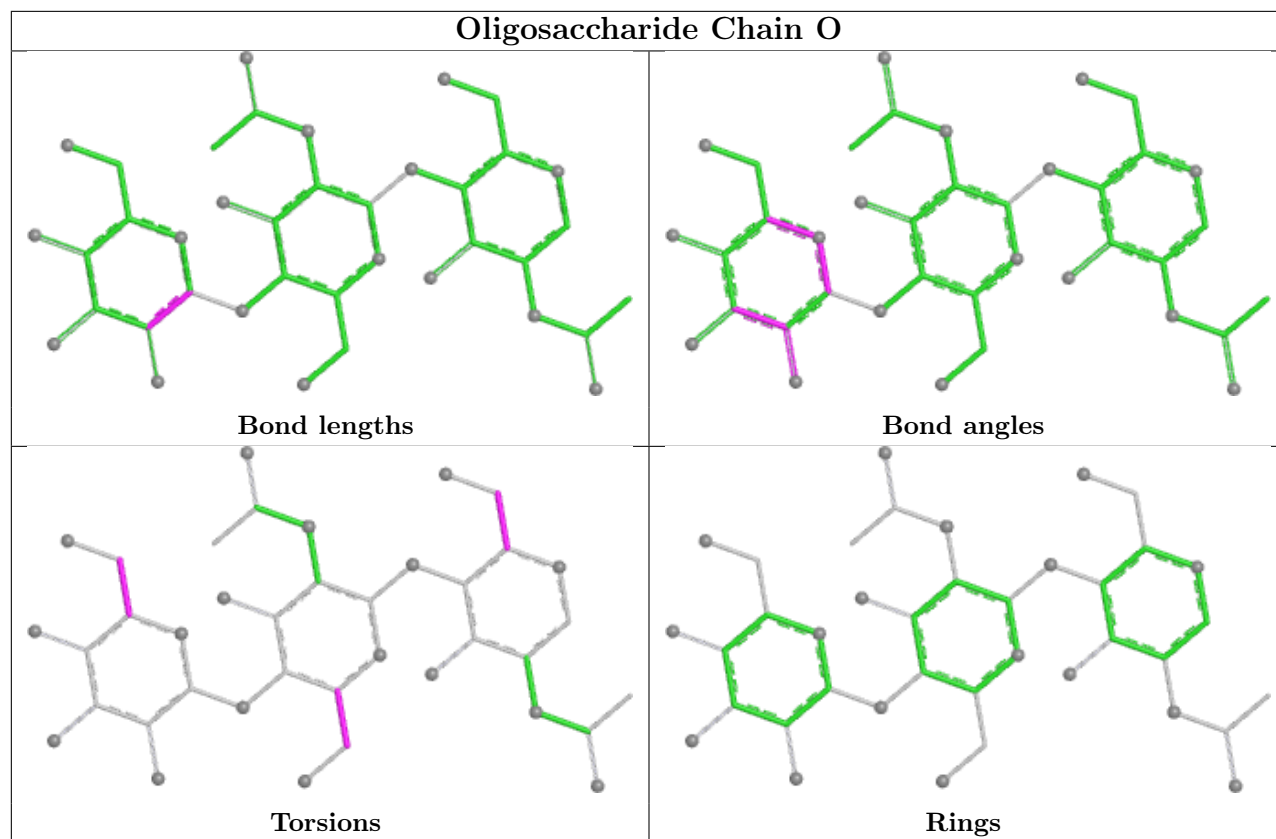


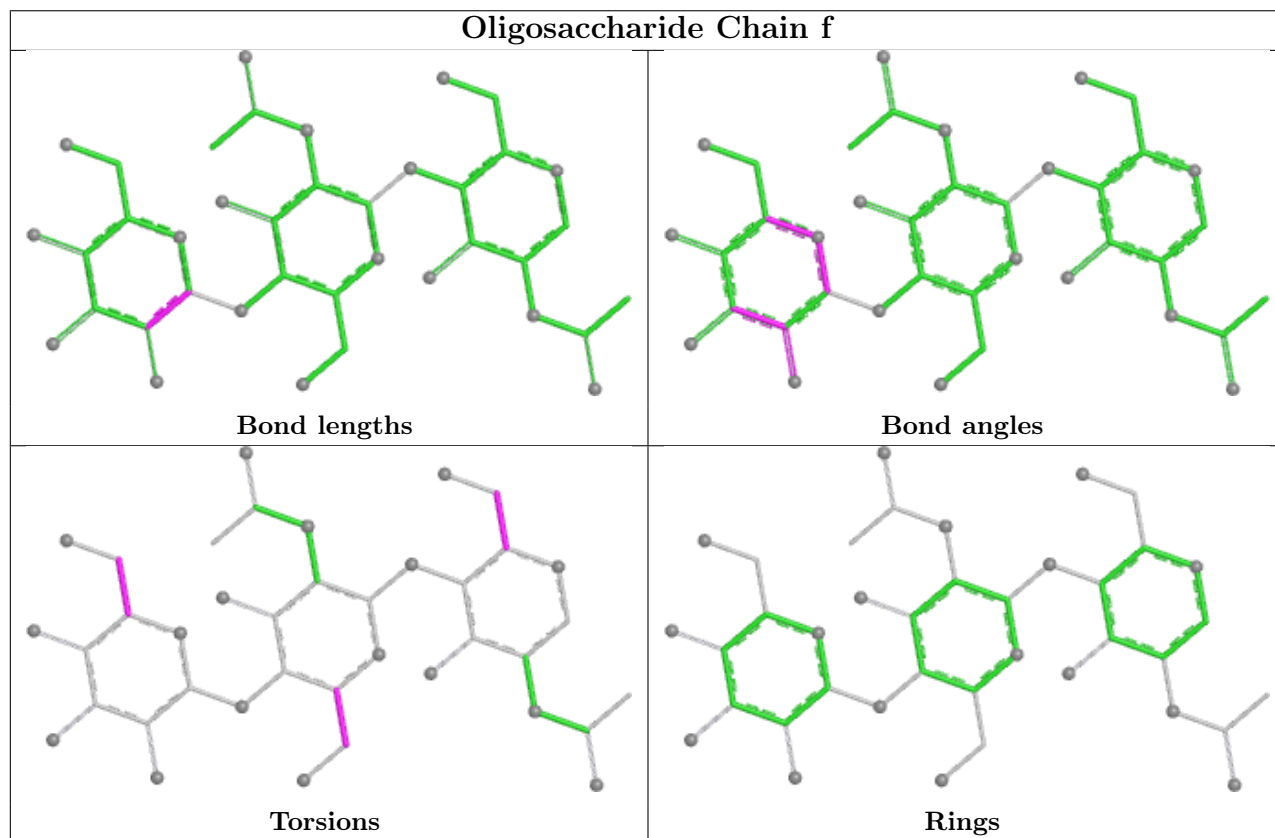
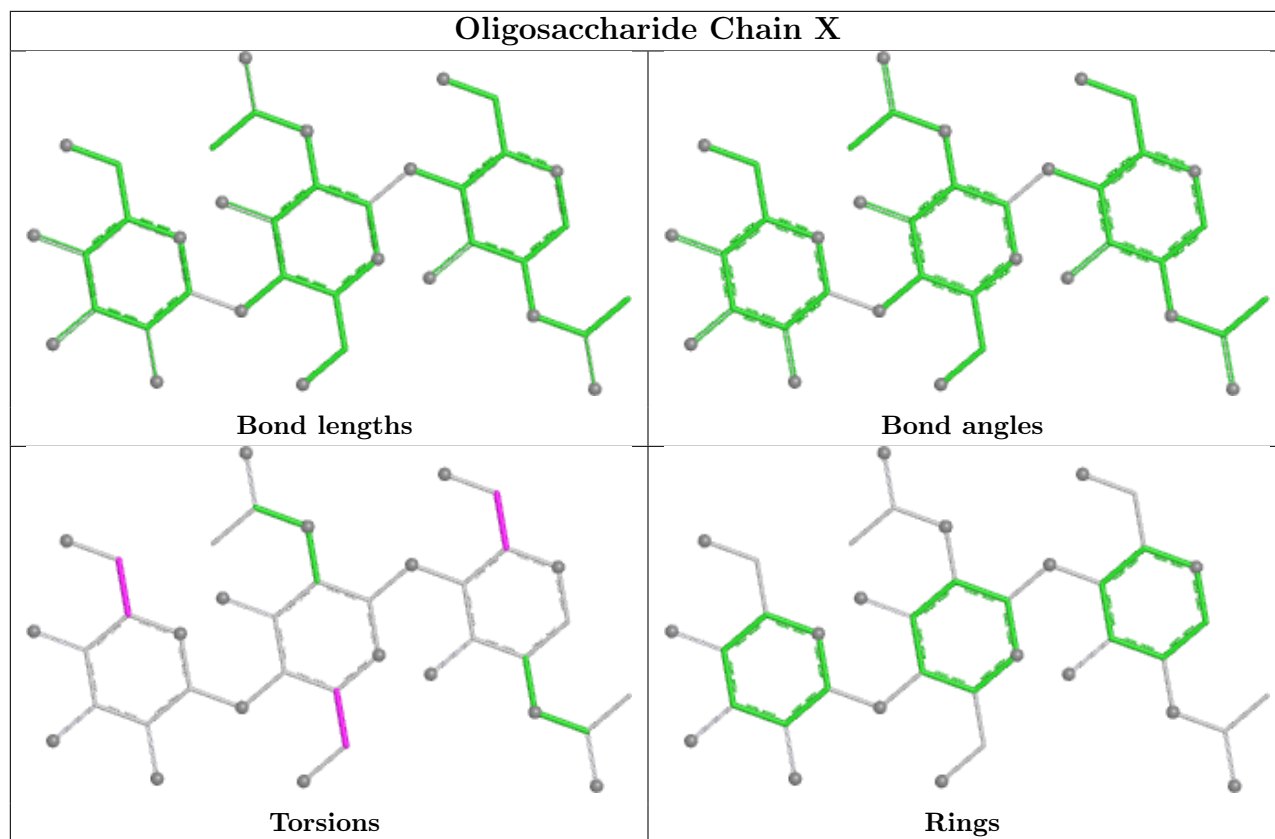


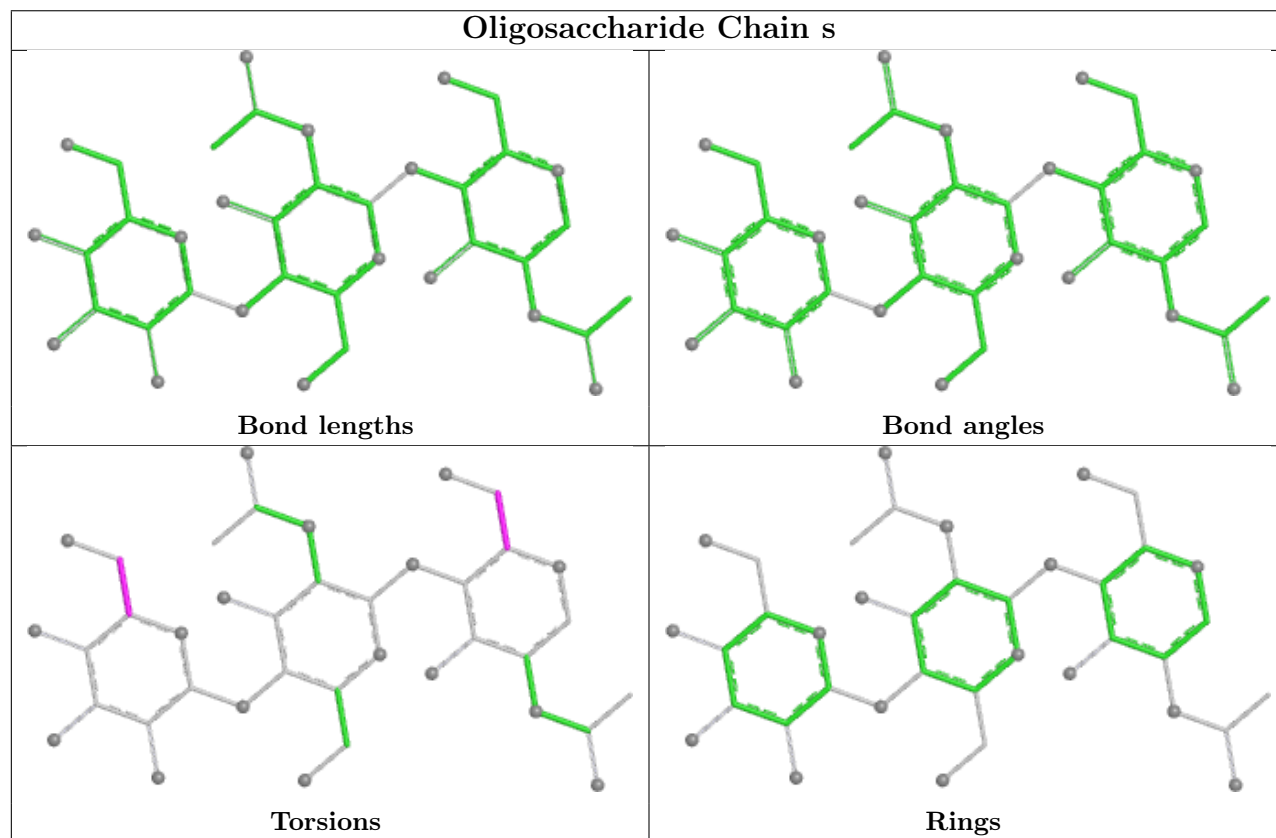
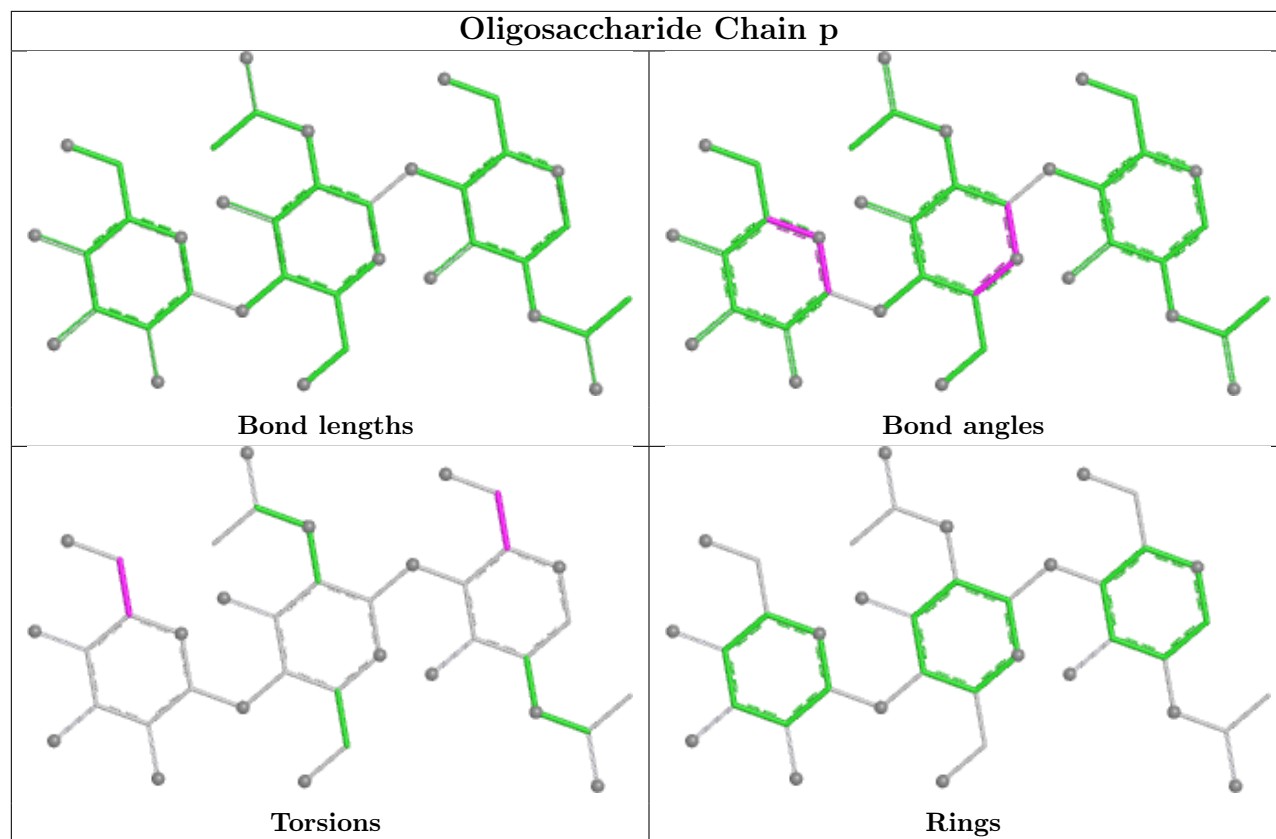


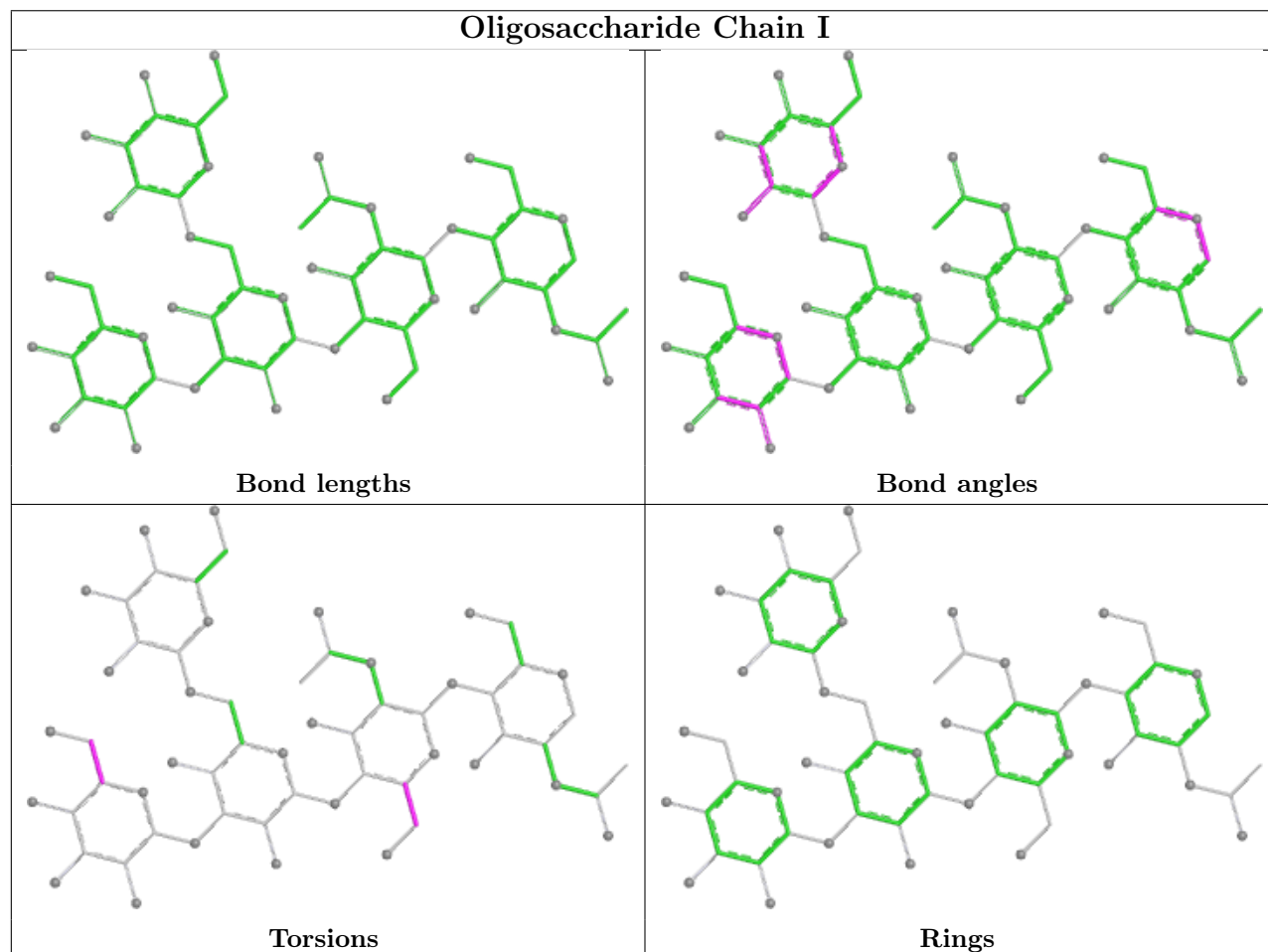
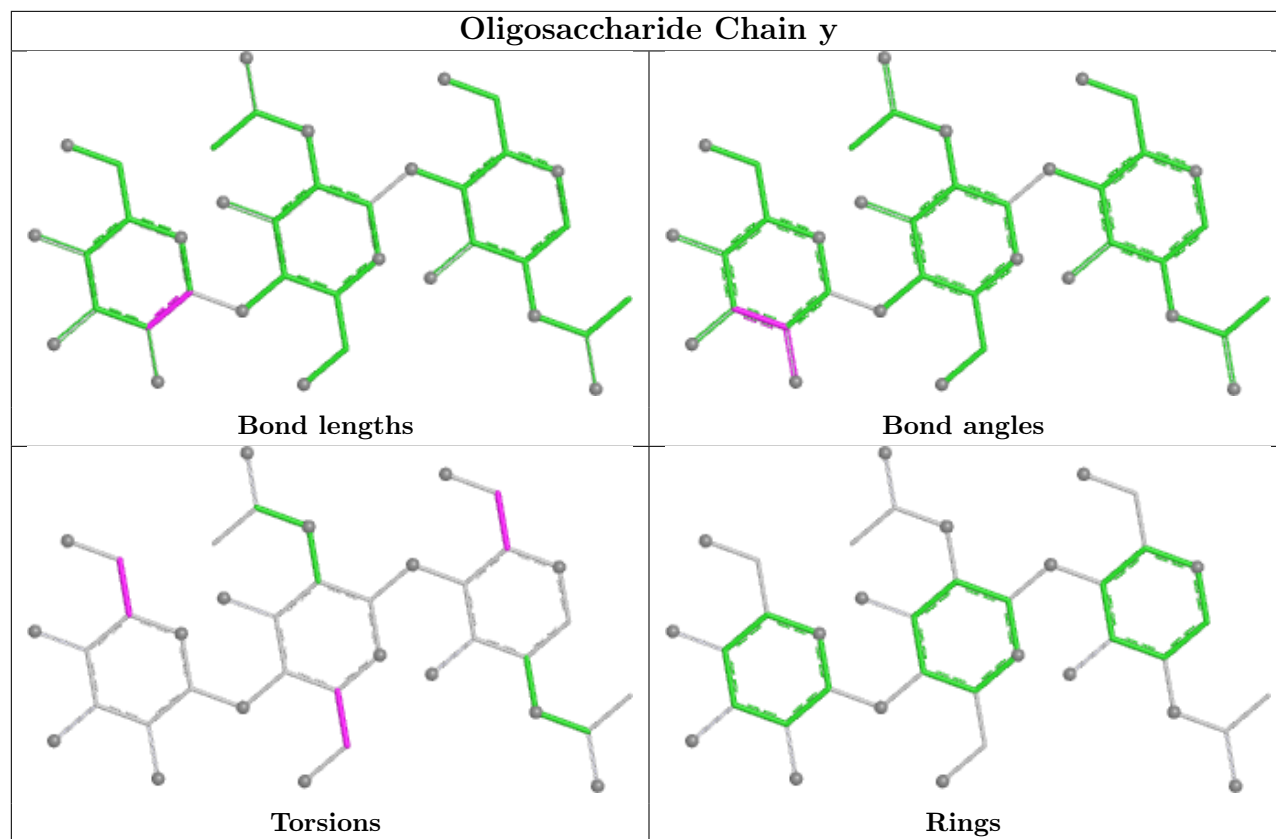


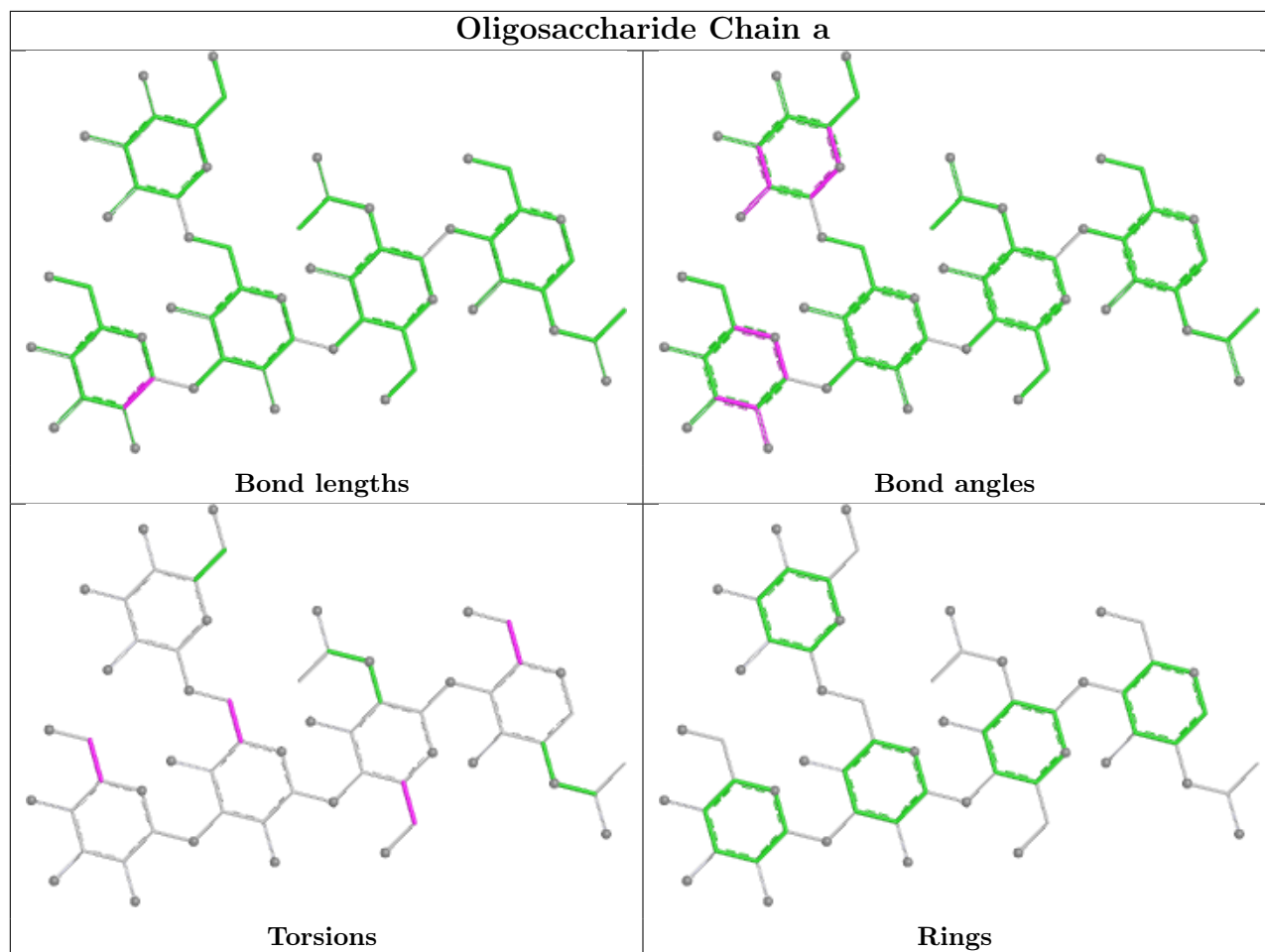


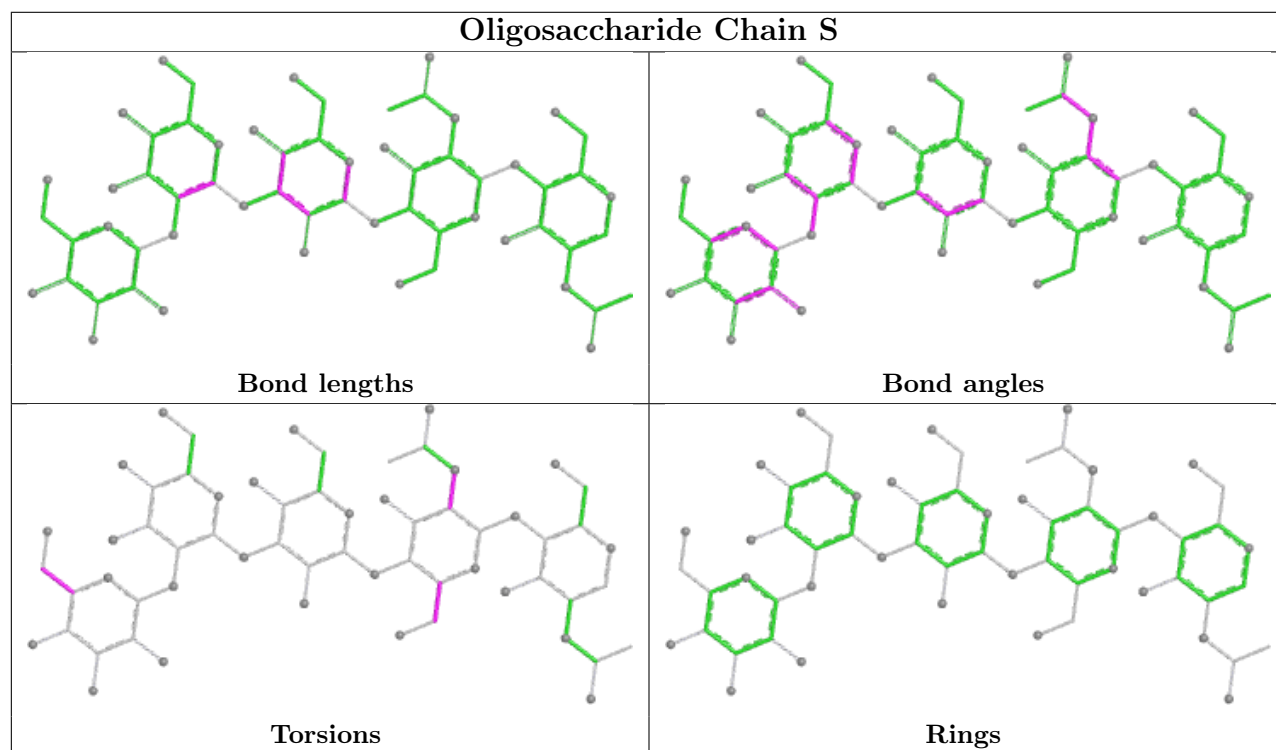
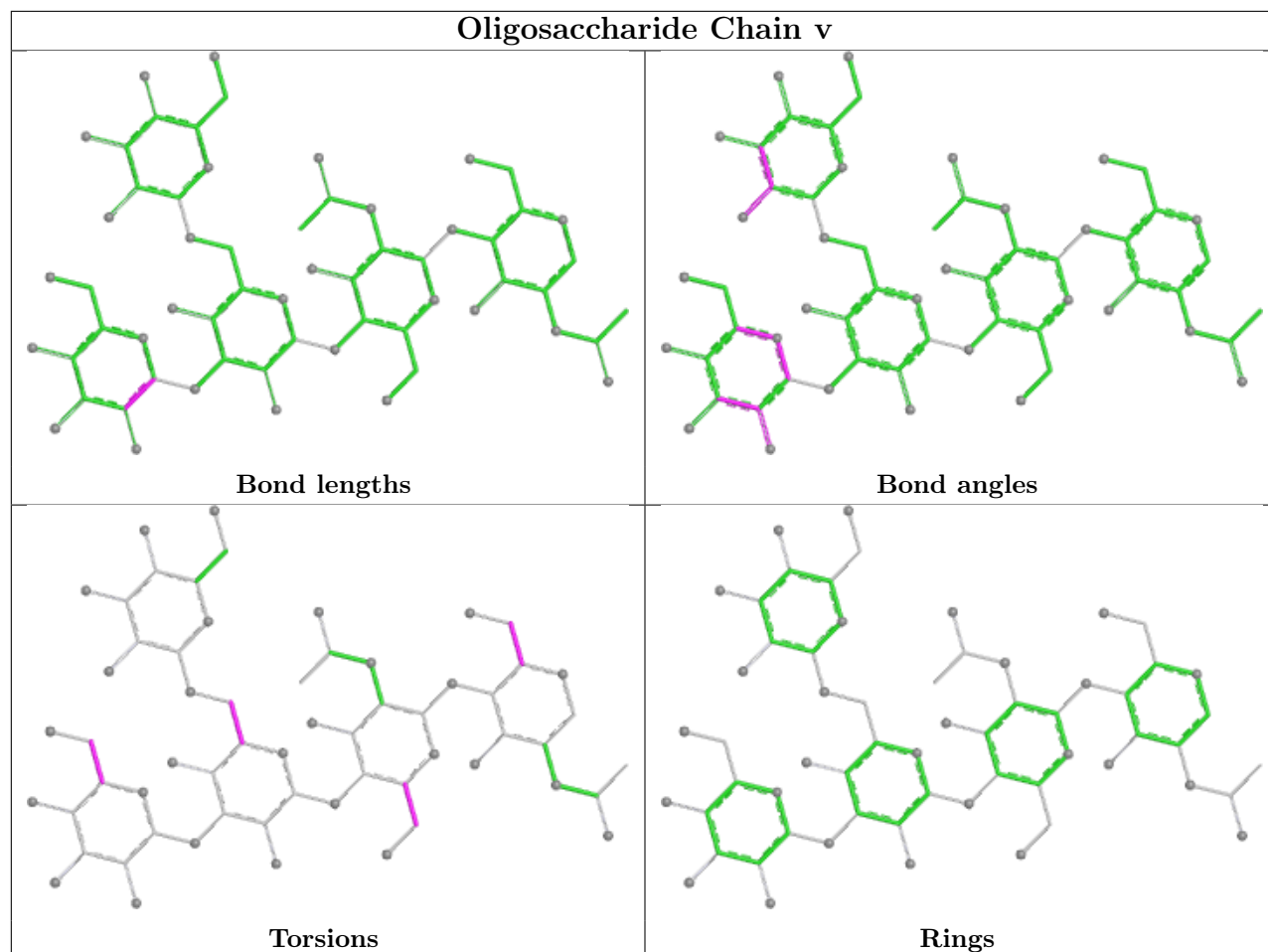


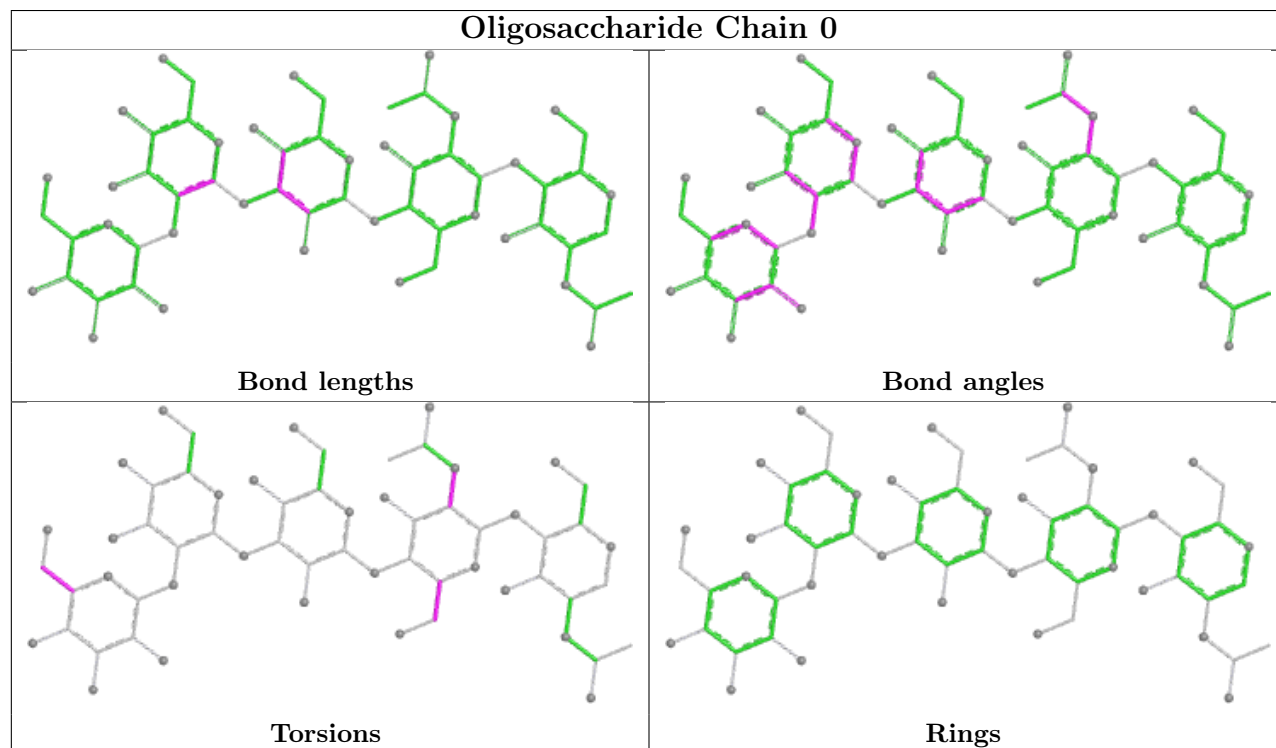
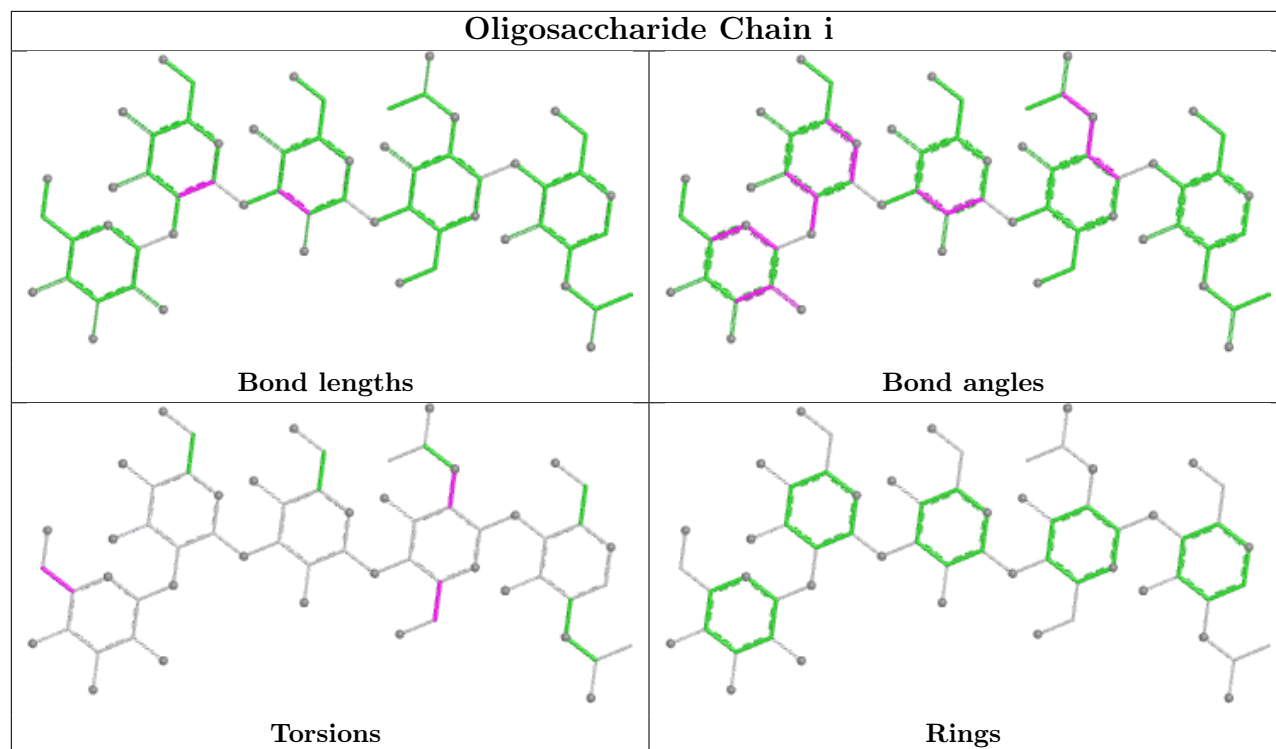


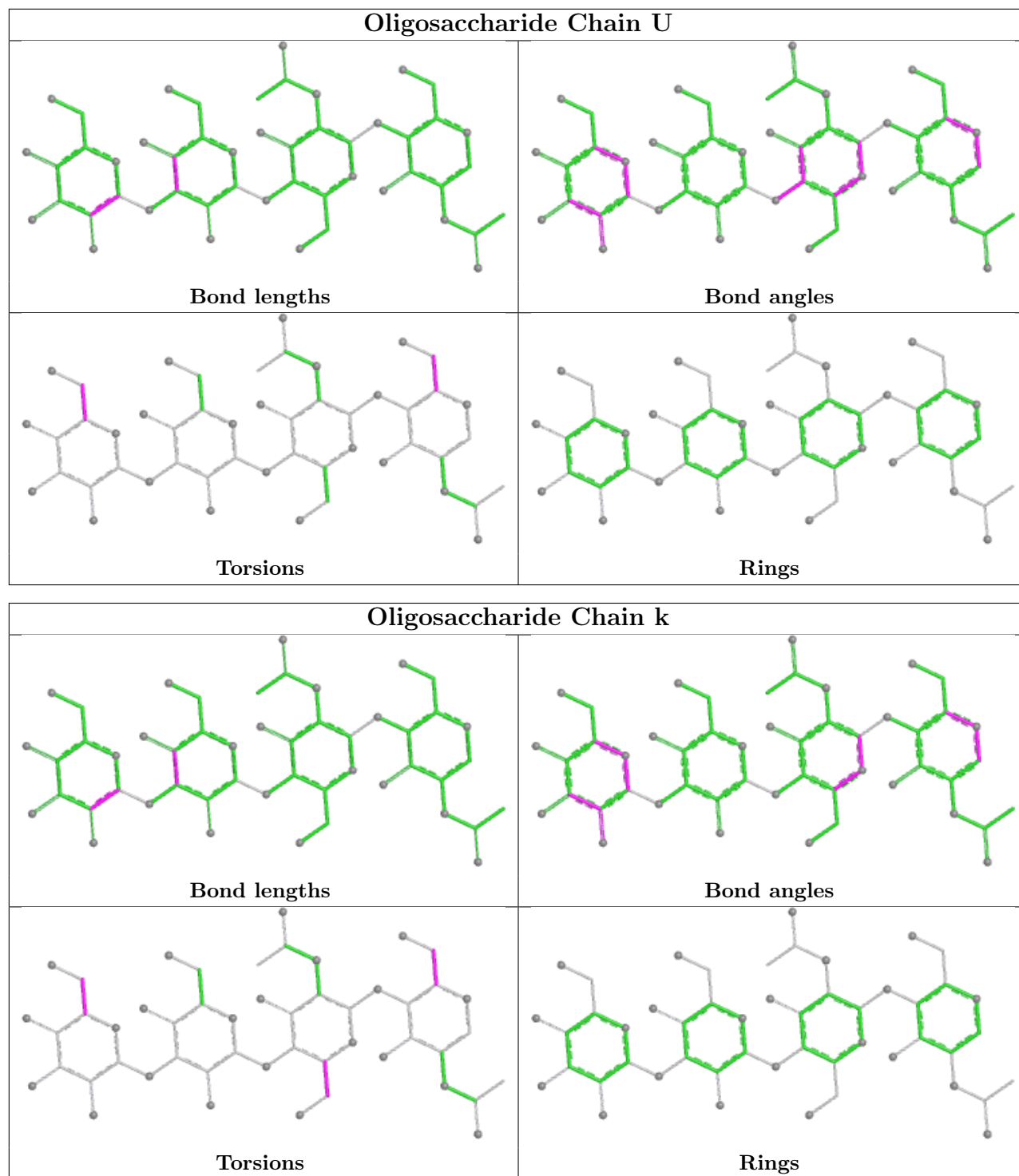


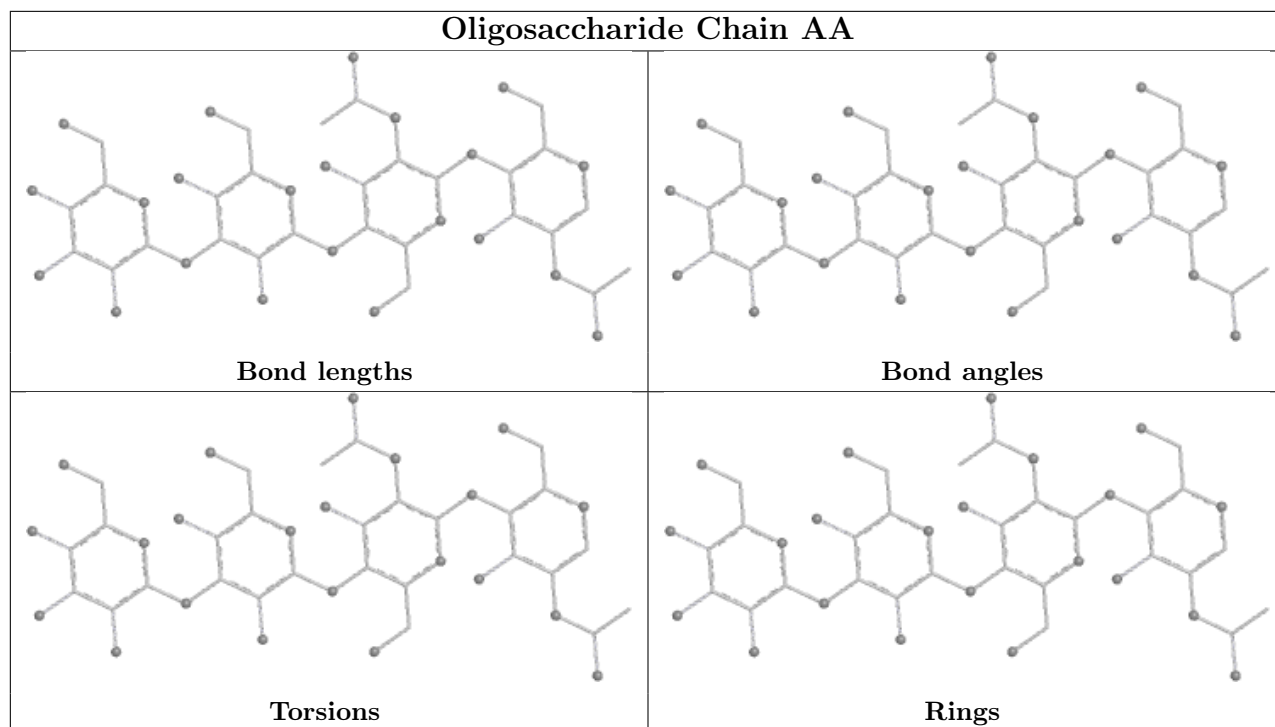












5.6 Ligand geometry [i](#)

3 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
14	NAG	c	611	1	14,14,15	0.24	0	17,19,21	0.60	1 (5%)
14	NAG	2	611	1	14,14,15	0.22	0	17,19,21	0.57	0
14	NAG	C	611	1	14,14,15	0.22	0	17,19,21	0.55	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
14	NAG	c	611	1	-	0/6/23/26	0/1/1/1
14	NAG	2	611	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	NAG	C	611	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	c	611	NAG	C1-O5-C5	2.01	114.89	112.19

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
14	C	611	NAG	C4-C5-C6-O6
14	C	611	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

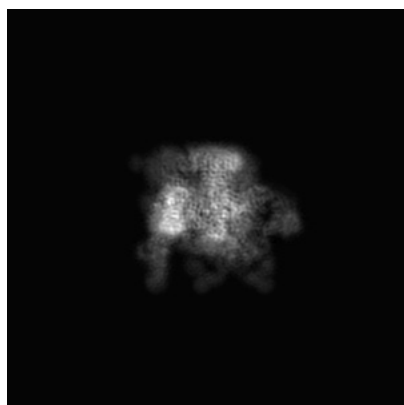
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-7621. These allow visual inspection of the internal detail of the map and identification of artifacts.

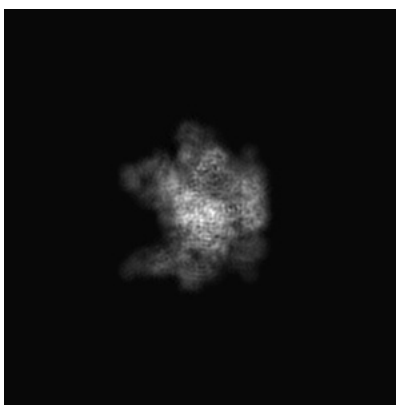
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

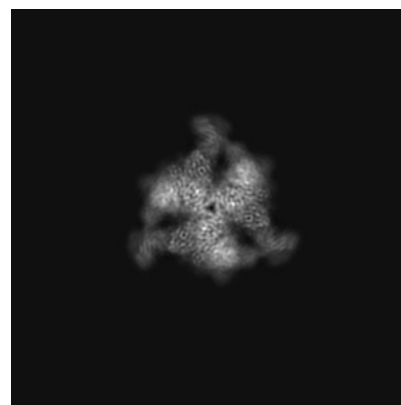
6.1.1 Primary map



X



Y

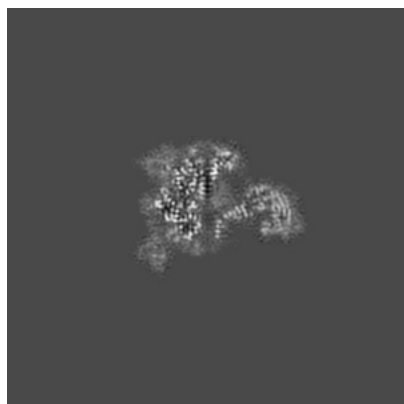


Z

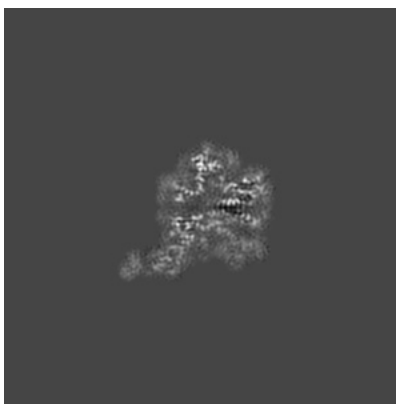
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 192



Y Index: 192

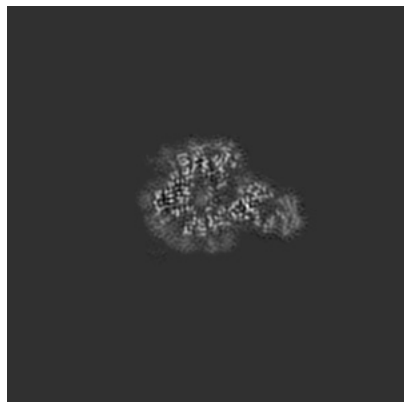


Z Index: 192

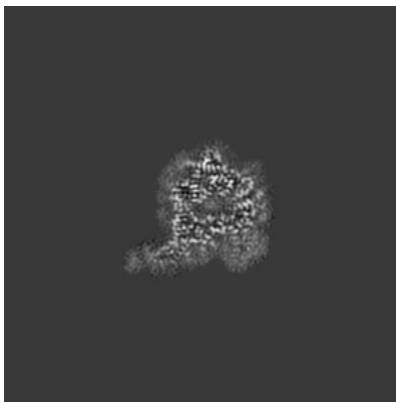
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

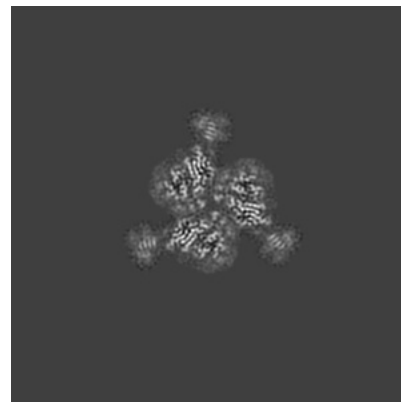
6.3.1 Primary map



X Index: 184



Y Index: 201

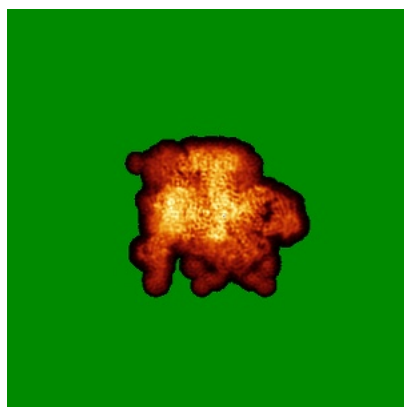


Z Index: 189

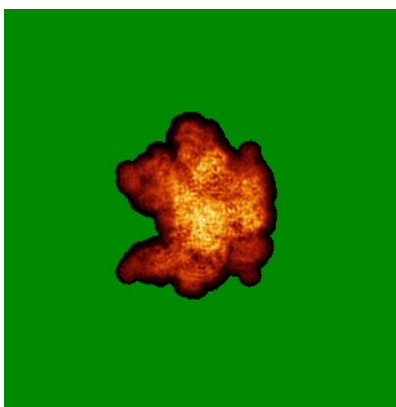
The images above show the largest variance slices of the map in three orthogonal directions.

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

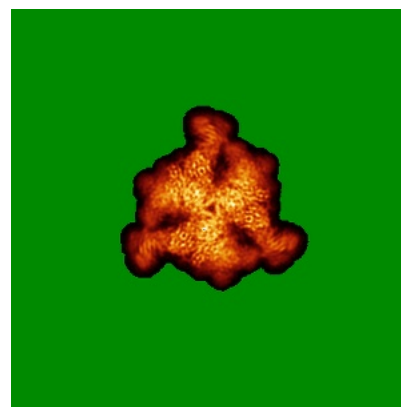
6.4.1 Primary map



X



Y

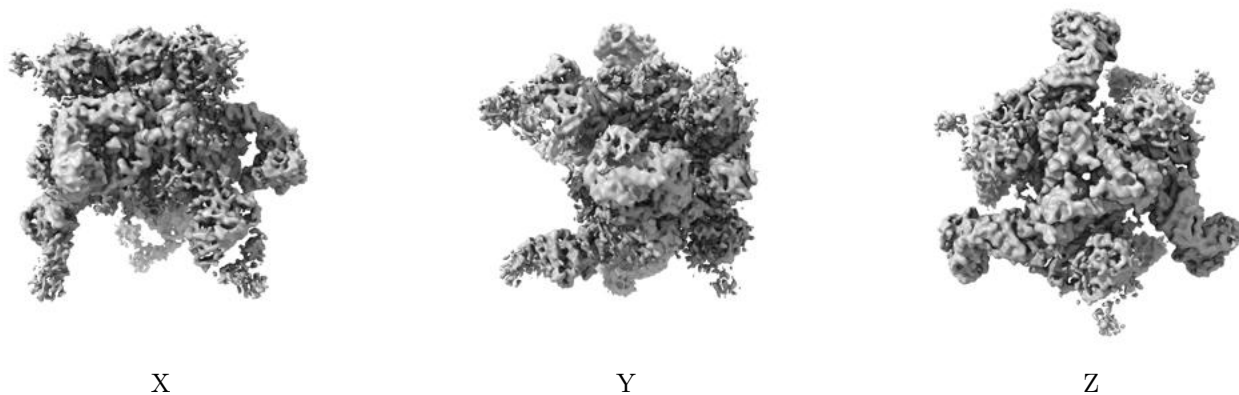


Z

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 1.04. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

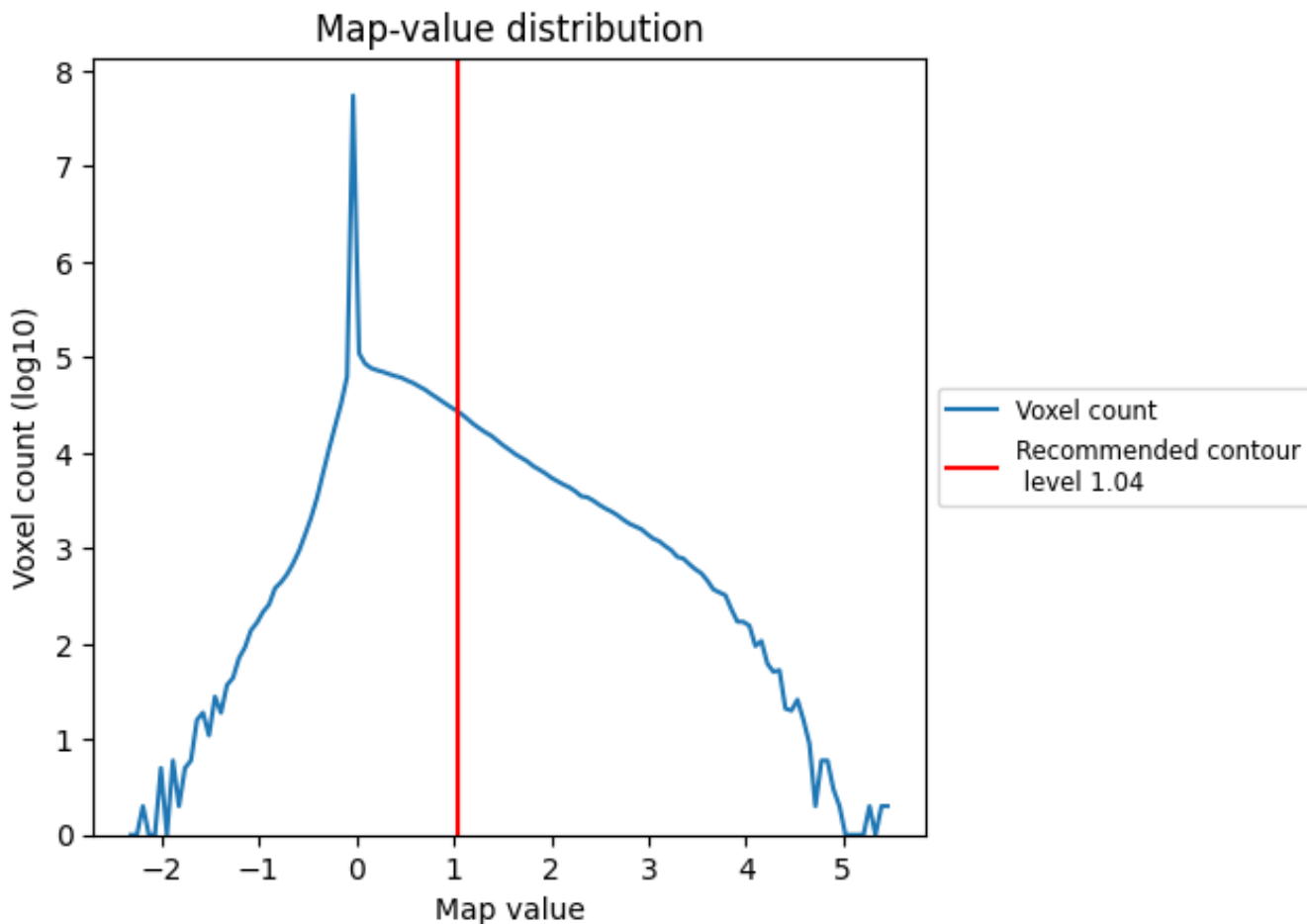
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

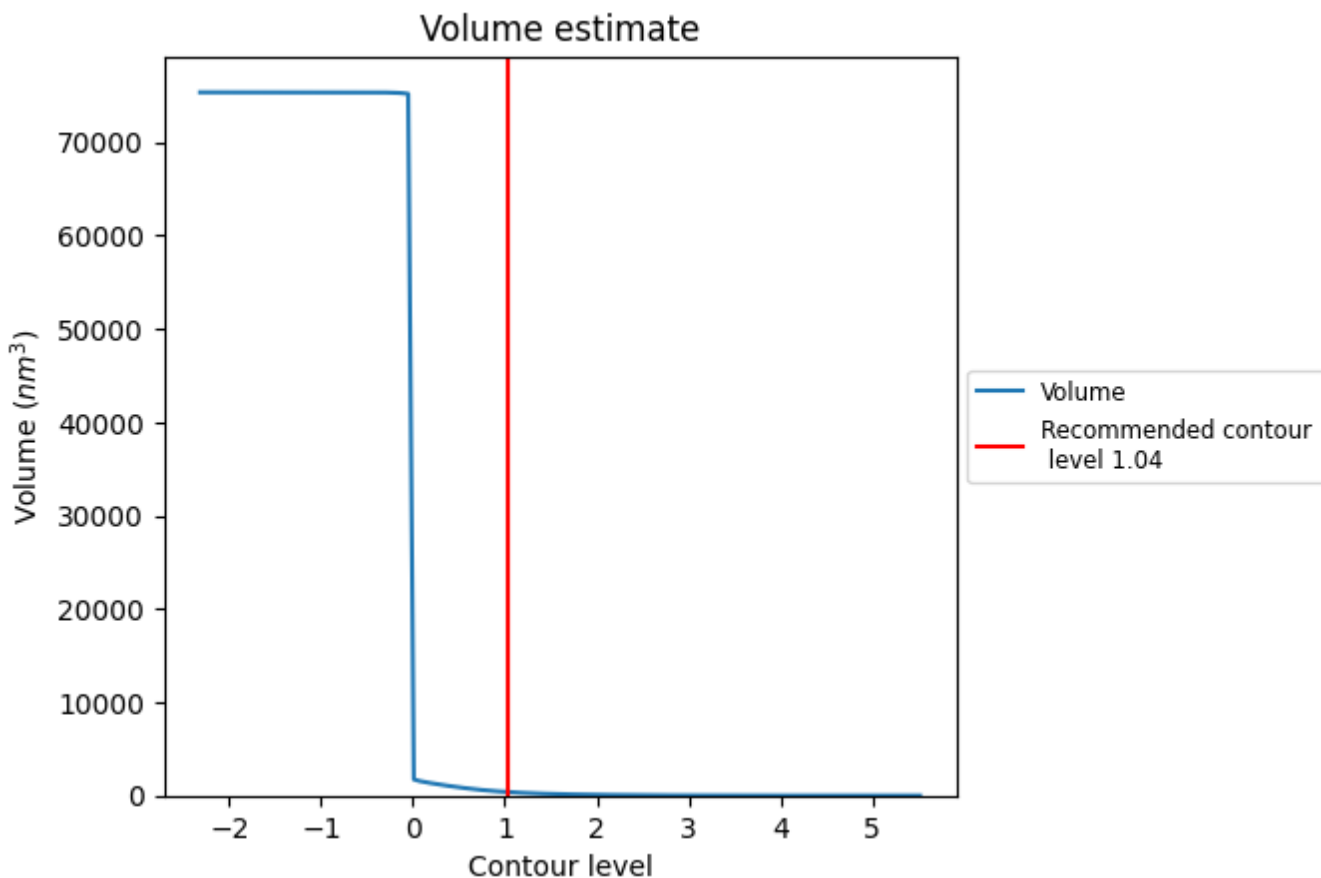
This section contains the results of statistical analysis of the map.

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.

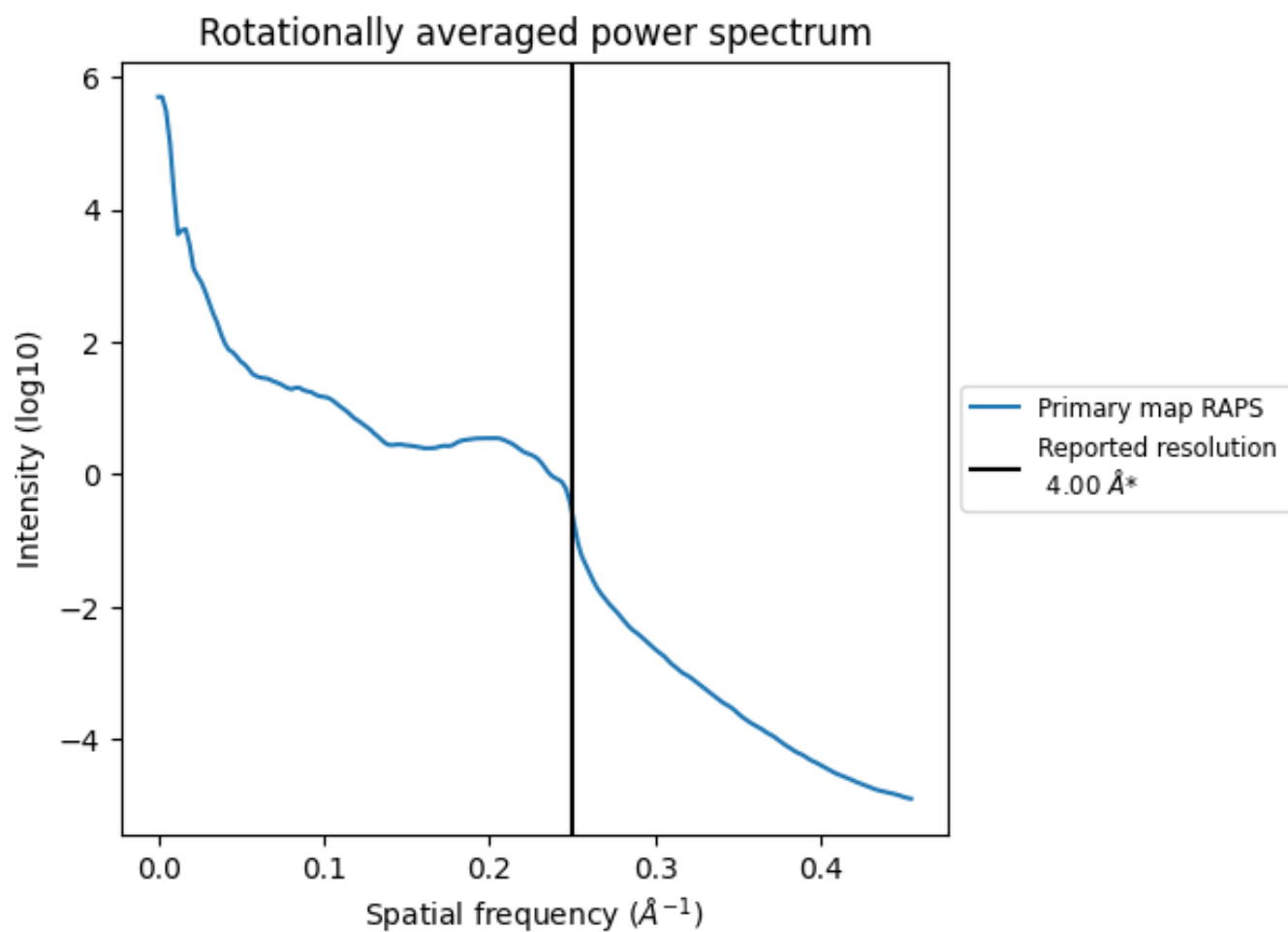
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 374 nm³; this corresponds to an approximate mass of 338 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 [i](#)



*Reported resolution corresponds to spatial frequency of 0.250\AA^{-1}

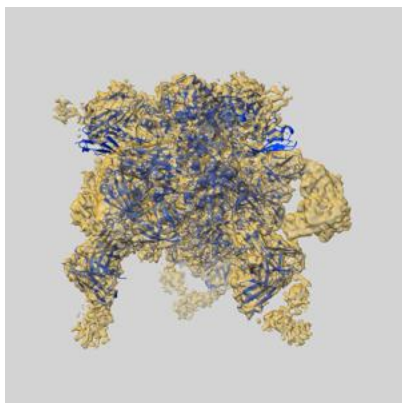
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

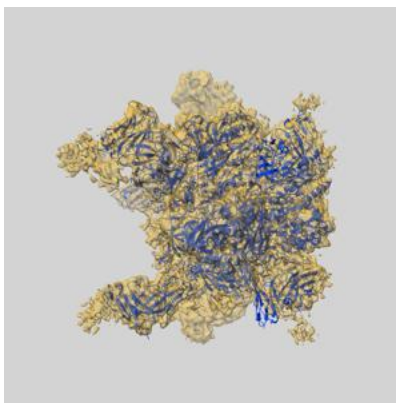
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-7621 and PDB model 6CUE. Per-residue inclusion information can be found in section 3 on page 13.

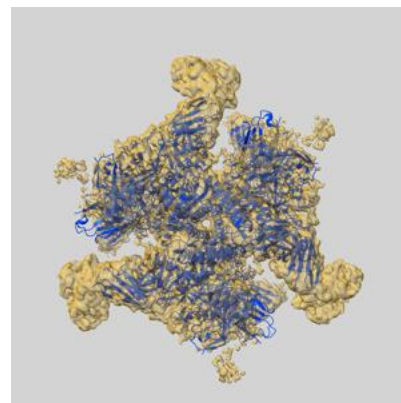
9.1 Map-model overlay [i](#)



X



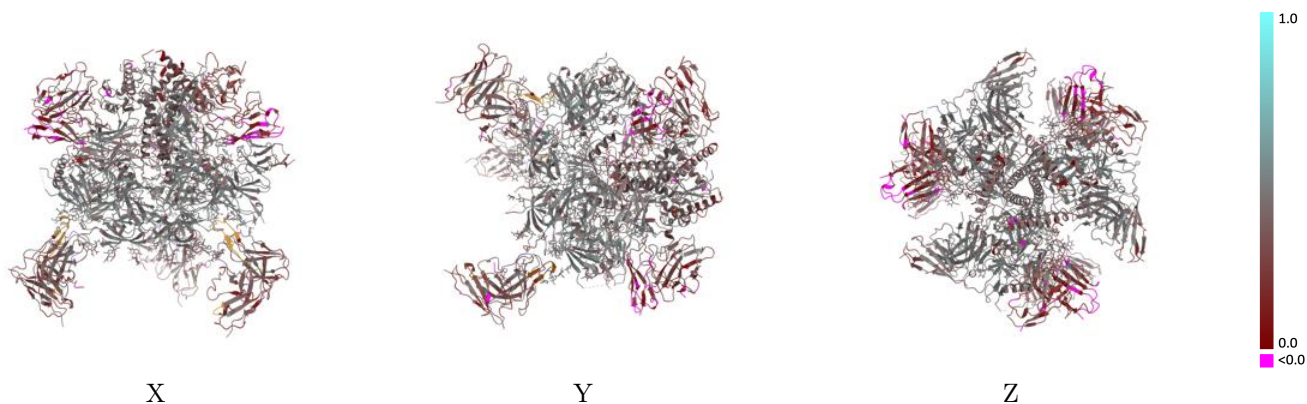
Y



Z

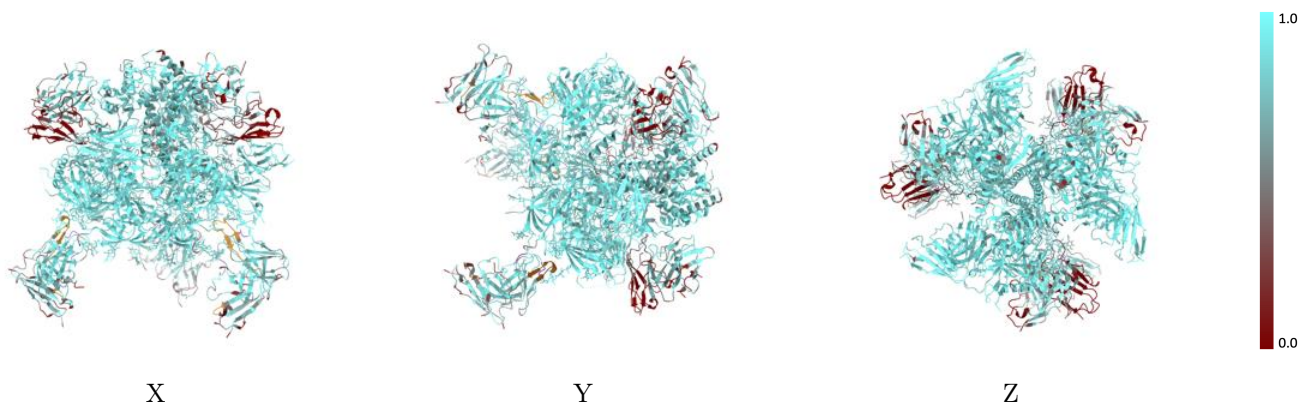
The images above show the 3D surface view of the map at the recommended contour level 1.04 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



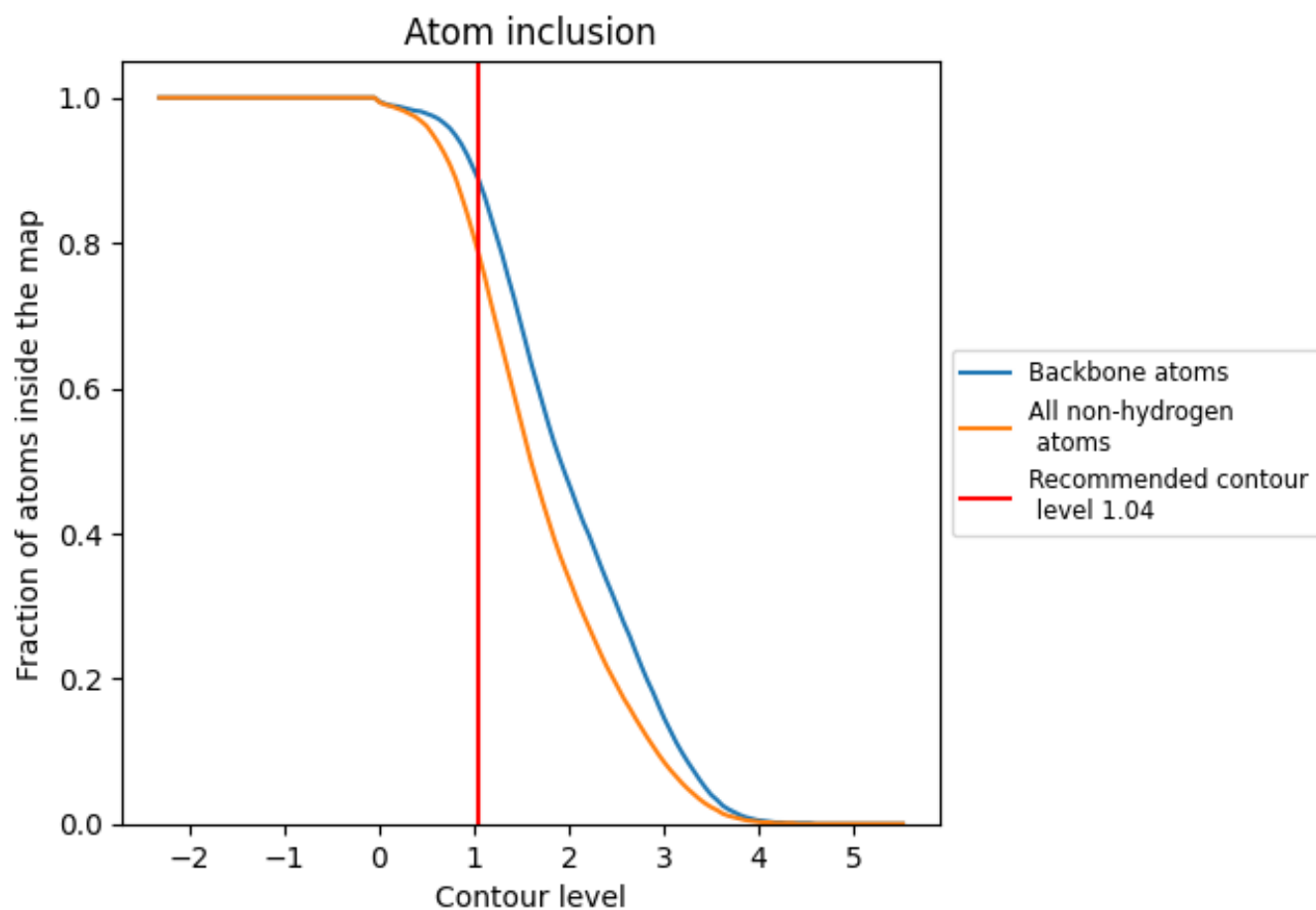
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (1.04).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 89% of all backbone atoms, 79% 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 (1.04) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7890	 0.3790
0	 0.9840	 0.4710
1	 0.8520	 0.3820
2	 0.8950	 0.4460
3	 0.6360	 0.2630
4	 0.2940	 0.1720
5	 0.7120	 0.3140
6	 0.7390	 0.3450
7	 0.9050	 0.4600
8	 0.8930	 0.4120
9	 0.6790	 0.2950
A	 0.7860	 0.3410
AA	 0.8400	 0.4550
B	 0.7440	 0.4020
C	 0.8930	 0.4420
D	 0.8590	 0.3760
E	 0.9230	 0.4560
F	 0.7500	 0.4130
G	 0.9290	 0.4670
H	 0.6300	 0.2690
I	 0.8530	 0.4440
J	 0.8210	 0.4680
K	 0.7140	 0.3940
L	 0.2840	 0.1770
M	 0.7160	 0.3330
N	 0.7660	 0.3530
O	 0.8970	 0.4330
P	 0.8930	 0.4590
Q	 0.9090	 0.4600
R	 0.8900	 0.4050
S	 0.9670	 0.4810
T	 0.7140	 0.3650
U	 0.8200	 0.4590
V	 0.7500	 0.3570
W	 0.7690	 0.4270



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Chain	Atom inclusion	Q-score
X	 0.8970	 0.4580
Y	 0.8210	 0.3910
Z	 0.8930	 0.4660
a	 0.8690	 0.4540
b	 0.8570	 0.4850
c	 0.8920	 0.4420
d	 0.8590	 0.3800
e	 0.6790	 0.4220
f	 0.9230	 0.4460
g	 0.9640	 0.4800
h	 0.6260	 0.2590
i	 0.9840	 0.4800
j	 0.6790	 0.3430
k	 0.8200	 0.4530
l	 0.2860	 0.1580
m	 0.7020	 0.3250
n	 0.7640	 0.3500
o	 0.7860	 0.3610
p	 0.6920	 0.4100
q	 0.9110	 0.4590
r	 0.8900	 0.4060
s	 0.9230	 0.4660
t	 0.8210	 0.3930
u	 0.9290	 0.4660
v	 0.8690	 0.4380
w	 0.8570	 0.4610
x	 0.6790	 0.4160
y	 0.9490	 0.4580
z	 0.9290	 0.4430