



Full wwPDB EM Validation Report ⓘ

Mar 27, 2026 – 06:06 AM UTC

PDB ID : 6N1W / pdb_00006n1w
EMDB ID : EMD-9320
Title : Cryo-EM structure at 4.2 Å resolution of vaccine-elicited antibody DFPH-a.15 in complex with HIV-1 Env BG505 DS-SOSIP, and antibodies VRC03 and PGT122
Authors : Acharya, P.; Kwong, P.D.
Deposited on : 2018-11-12
Resolution : 4.20 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:

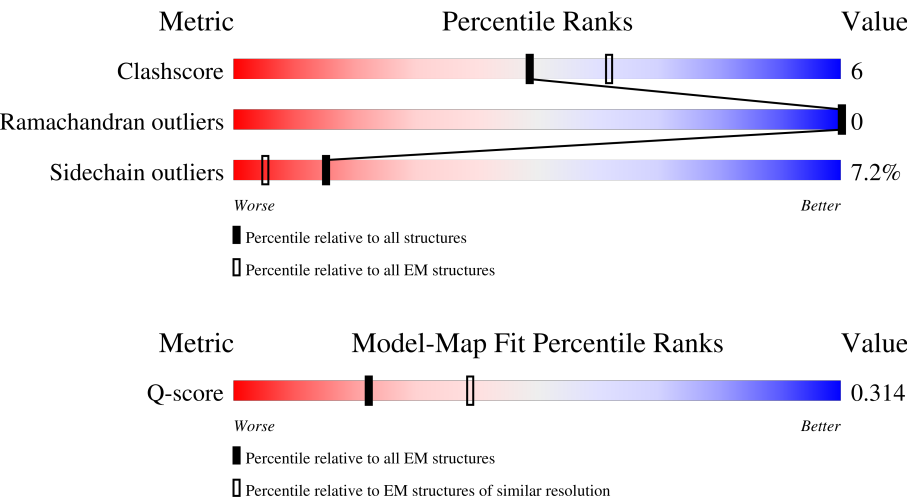
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 4.20 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	5410 (3.70 - 4.70)

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	75% 19% ..
1	C	473	75% 19% ..
1	c	473	75% 19% ..
2	3	224	17% 41% 15% 43% .

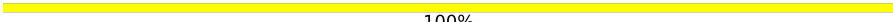
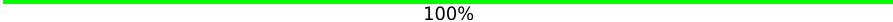
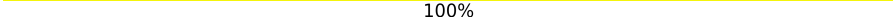
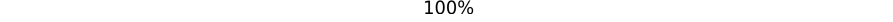
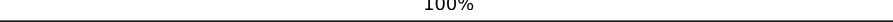
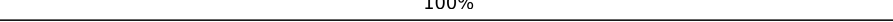
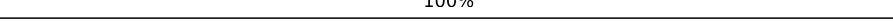
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Mol	Chain	Length	Quality of chain
2	H	224	
2	h	224	
3	4	212	
3	L	212	
3	l	212	
4	5	132	
4	M	132	
4	m	132	
5	6	105	
5	N	105	
5	n	105	
6	7	102	
6	R	102	
6	r	102	
7	8	128	
7	Q	128	
7	q	128	
8	A	153	
8	D	153	
8	d	153	
9	9	2	
9	B	2	
9	F	2	
9	G	2	
9	I	2	

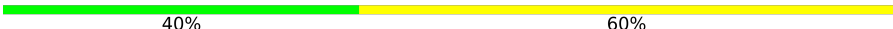
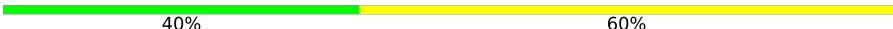
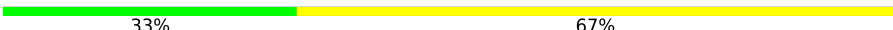
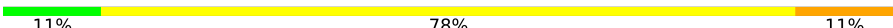
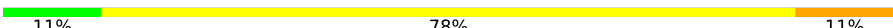
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Mol	Chain	Length	Quality of chain
9	K	2	 100%
9	O	2	 50% 50%
9	U	2	 100%
9	W	2	 50% 50%
9	Y	2	 100%
9	Z	2	 50% 50%
9	a	2	 100%
9	e	2	 100%
9	f	2	 50% 50%
9	k	2	 100%
9	p	2	 50% 50%
9	t	2	 100%
9	u	2	 50% 50%
9	v	2	 100%
9	x	2	 100%
9	y	2	 100%
10	0	4	 25% 75%
10	AA	4	 50% 25% 25%
10	E	4	 25% 25% 75%
10	S	4	 25% 75%
10	V	4	 25% 50% 25%
10	X	4	 25% 75%
10	i	4	 100%
10	o	4	 50% 25% 25%
10	s	4	 25% 25% 75%

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Mol	Chain	Length	Quality of chain
11	J	5	 40%60%
11	b	5	 40%60%
11	w	5	 40%60%
12	P	3	 67%33%
12	g	3	 33%67%
12	z	3	 67%33%
13	1	9	 11%89%
13	T	9	 11%78%11%
13	j	9	 11%78%11%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 31905 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
1	2	453	Total	C	N	O	S	0	0
			3564	2233	630	671	30		
1	C	453	Total	C	N	O	S	0	0
			3564	2233	630	671	30		
1	c	453	Total	C	N	O	S	0	0
			3564	2233	630	671	30		

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
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
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 DFPH-a.15 heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	3	127	Total	C	N	O	S	0	0
			980	617	167	193	3		
2	H	127	Total	C	N	O	S	0	0
			980	617	167	193	3		
2	h	127	Total	C	N	O	S	0	0
			980	617	167	193	3		

- Molecule 3 is a protein called DFPH-a.15 Light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	4	105	Total	C	N	O	S	0	0
			787	491	130	163	3		
3	L	105	Total	C	N	O	S	0	0
			787	491	130	163	3		
3	l	105	Total	C	N	O	S	0	0
			787	491	130	163	3		

- Molecule 4 is a protein called PGT122 Heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	5	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		
4	M	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		
4	m	132	Total	C	N	O	S	0	0
			1047	669	180	195	3		

- Molecule 5 is a protein called PGT122 Light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	6	105	Total	C	N	O	S	0	0
			805	504	139	160	2		
5	N	105	Total	C	N	O	S	0	0
			805	504	139	160	2		
5	n	105	Total	C	N	O	S	0	0
			805	504	139	160	2		

- Molecule 6 is a protein called VRC03 Light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	7	102	Total	C	N	O	S	0	0
			802	510	137	152	3		
6	R	102	Total	C	N	O	S	0	0
			802	510	137	152	3		
6	r	102	Total	C	N	O	S	0	0
			802	510	137	152	3		

- Molecule 7 is a protein called VRC03 Heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	8	128	Total	C	N	O	S	0	0
			1023	657	175	185	6		
7	Q	128	Total	C	N	O	S	0	0
			1023	657	175	185	6		
7	q	128	Total	C	N	O	S	0	0
			1023	657	175	185	6		

- Molecule 8 is a protein called Envelope glycoprotein gp41.

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

There are 3 discrepancies between the modelled and reference sequences:

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

- 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	B	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	I	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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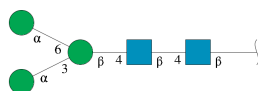
Mol	Chain	Residues	Atoms				AltConf	Trace
9	O	2	Total	C	N	O	0	0
			28	16	2	10		
9	U	2	Total	C	N	O	0	0
			28	16	2	10		
9	W	2	Total	C	N	O	0	0
			28	16	2	10		
9	Y	2	Total	C	N	O	0	0
			28	16	2	10		
9	Z	2	Total	C	N	O	0	0
			28	16	2	10		
9	a	2	Total	C	N	O	0	0
			28	16	2	10		
9	e	2	Total	C	N	O	0	0
			28	16	2	10		
9	f	2	Total	C	N	O	0	0
			28	16	2	10		
9	k	2	Total	C	N	O	0	0
			28	16	2	10		
9	p	2	Total	C	N	O	0	0
			28	16	2	10		
9	t	2	Total	C	N	O	0	0
			28	16	2	10		
9	u	2	Total	C	N	O	0	0
			28	16	2	10		
9	v	2	Total	C	N	O	0	0
			28	16	2	10		
9	x	2	Total	C	N	O	0	0
			28	16	2	10		
9	y	2	Total	C	N	O	0	0
			28	16	2	10		
9	9	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 10 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.



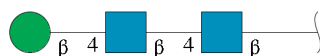
Mol	Chain	Residues	Atoms				AltConf	Trace
10	E	4	Total	C	N	O	0	0
			50	28	2	20		
10	S	4	Total	C	N	O	0	0
			50	28	2	20		
10	V	4	Total	C	N	O	0	0
			50	28	2	20		
10	X	4	Total	C	N	O	0	0
			50	28	2	20		
10	i	4	Total	C	N	O	0	0
			50	28	2	20		
10	o	4	Total	C	N	O	0	0
			50	28	2	20		
10	s	4	Total	C	N	O	0	0
			50	28	2	20		
10	0	4	Total	C	N	O	0	0
			50	28	2	20		
10	AA	4	Total	C	N	O	0	0
			50	28	2	20		

- 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.



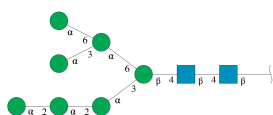
Mol	Chain	Residues	Atoms				AltConf	Trace
11	J	5	Total	C	N	O	0	0
			61	34	2	25		
11	b	5	Total	C	N	O	0	0
			61	34	2	25		
11	w	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 12 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.



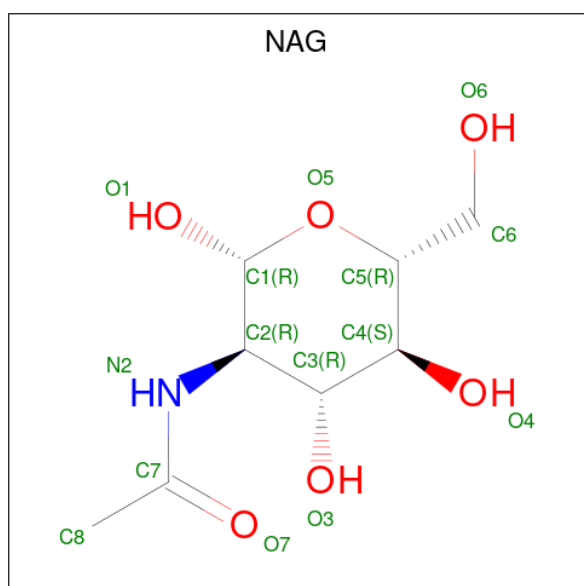
Mol	Chain	Residues	Atoms				AltConf	Trace
12	P	3	Total	C	N	O	0	0
			39	22	2	15		
12	g	3	Total	C	N	O	0	0
			39	22	2	15		
12	z	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 13 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
13	T	9	Total	C	N	O	0	0
			105	58	2	45		
13	j	9	Total	C	N	O	0	0
			105	58	2	45		
13	1	9	Total	C	N	O	0	0
			105	58	2	45		

- Molecule 14 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

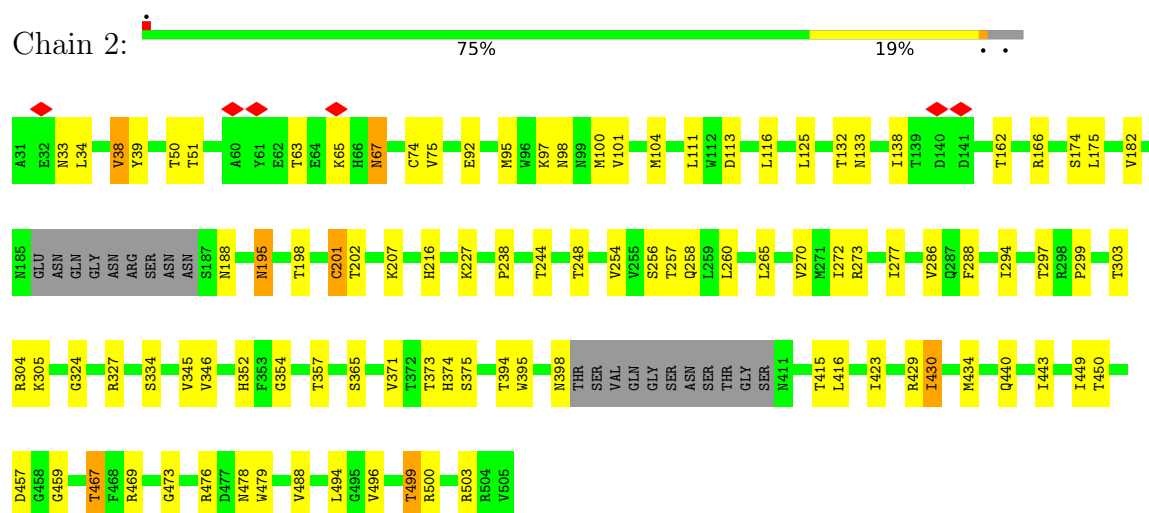


Mol	Chain	Residues	Atoms				AltConf
14	2	1	Total	C	N	O	0
			14	8	1	5	
14	2	1	Total	C	N	O	0
			14	8	1	5	
14	2	1	Total	C	N	O	0
			14	8	1	5	
14	C	1	Total	C	N	O	0
			14	8	1	5	
14	C	1	Total	C	N	O	0
			14	8	1	5	
14	C	1	Total	C	N	O	0
			14	8	1	5	
14	c	1	Total	C	N	O	0
			14	8	1	5	
14	c	1	Total	C	N	O	0
			14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

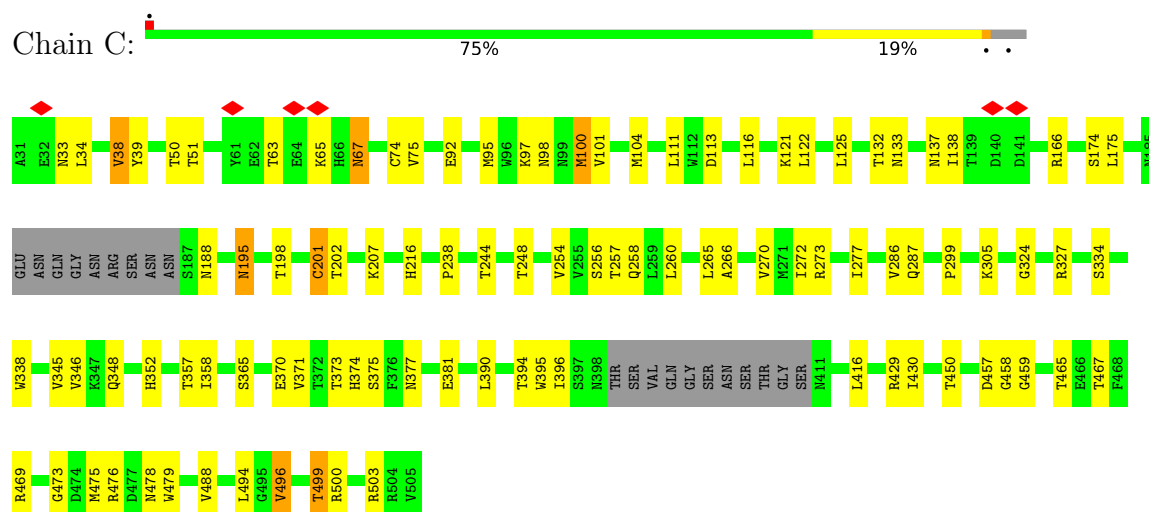
3 Residue-property plots [i](#)

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

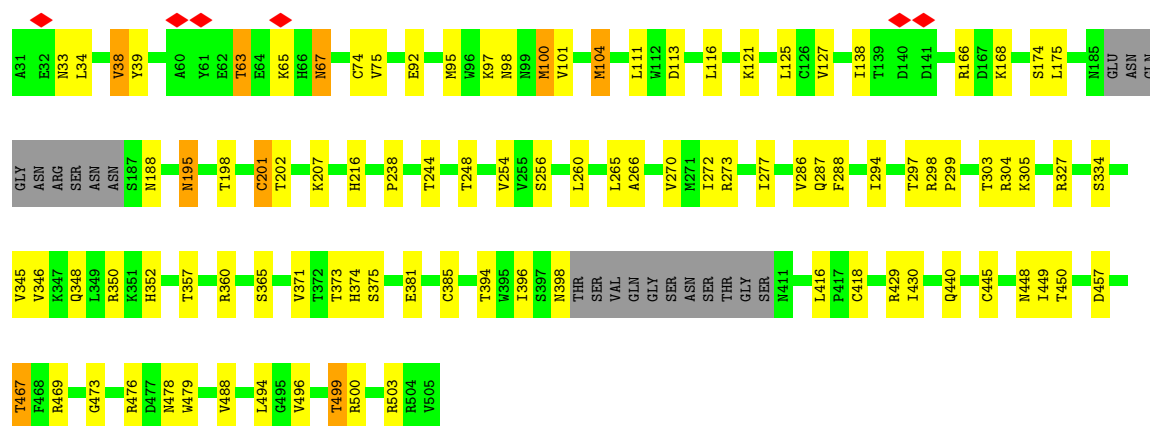


• Molecule 1: Envelope glycoprotein gp120

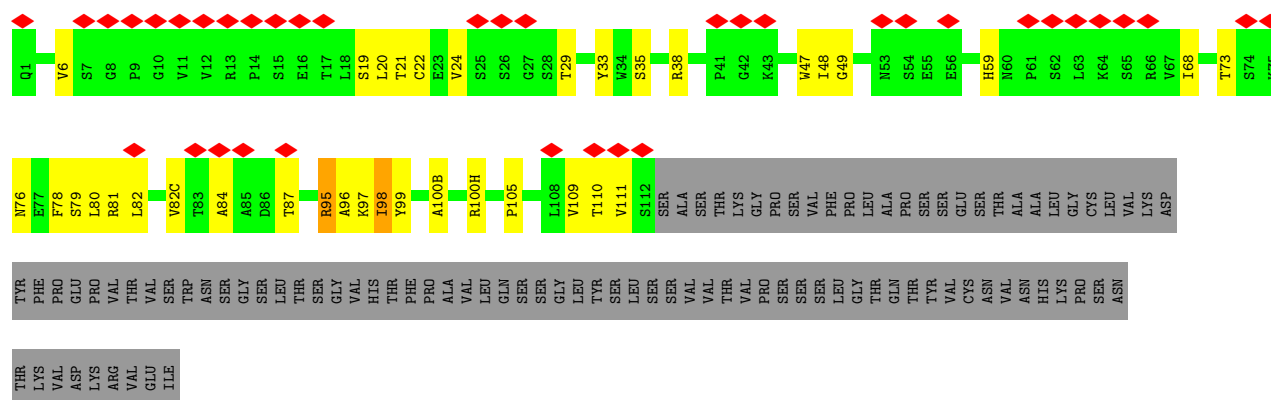


• Molecule 1: Envelope glycoprotein gp120

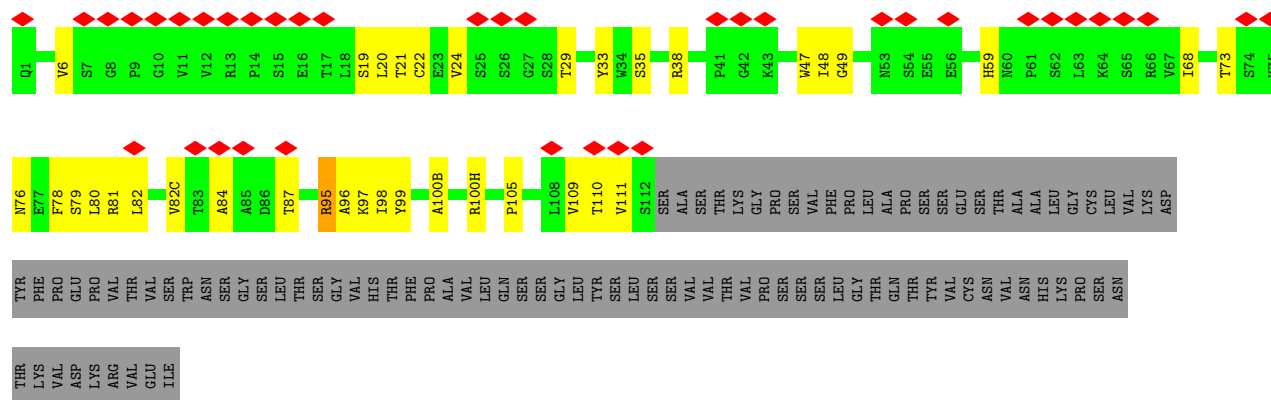
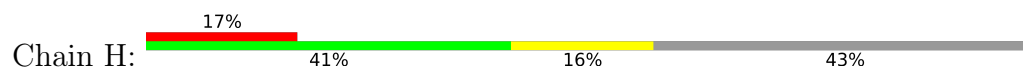




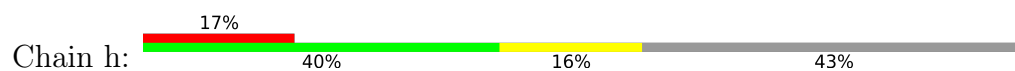
• Molecule 2: DFPH-a.15 heavy chain

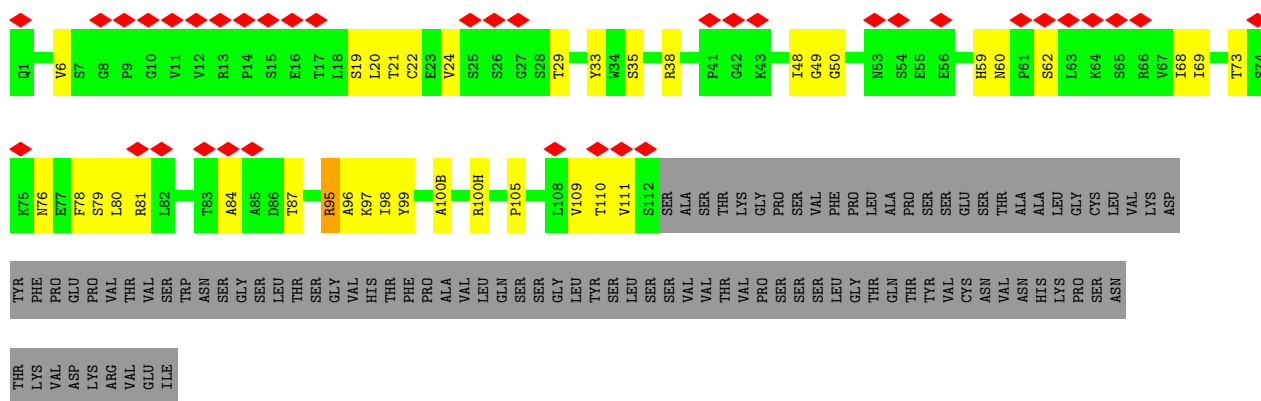


• Molecule 2: DFPH-a.15 heavy chain

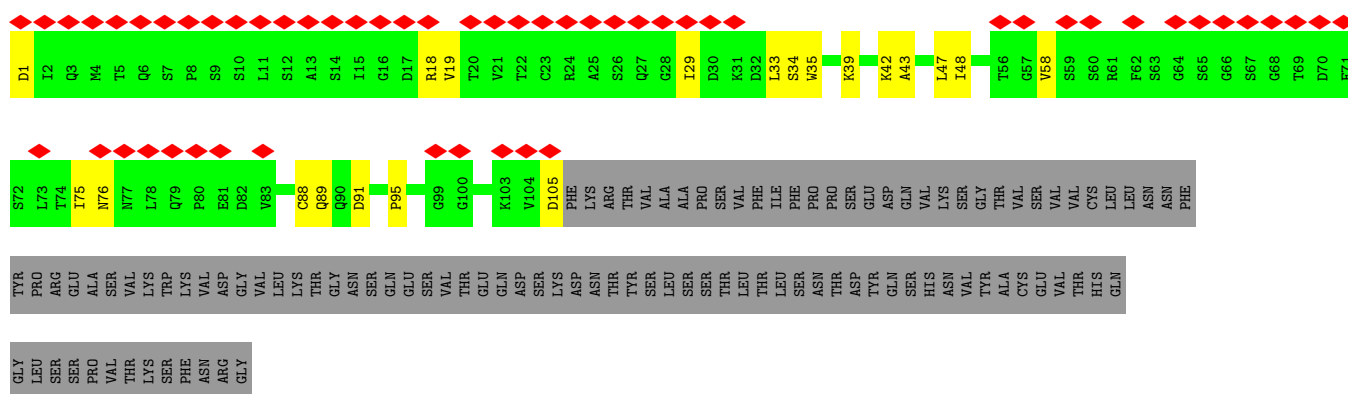
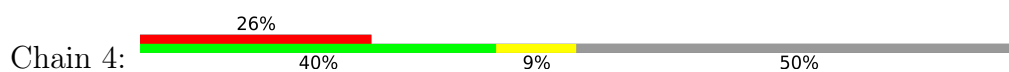


• Molecule 2: DFPH-a.15 heavy chain

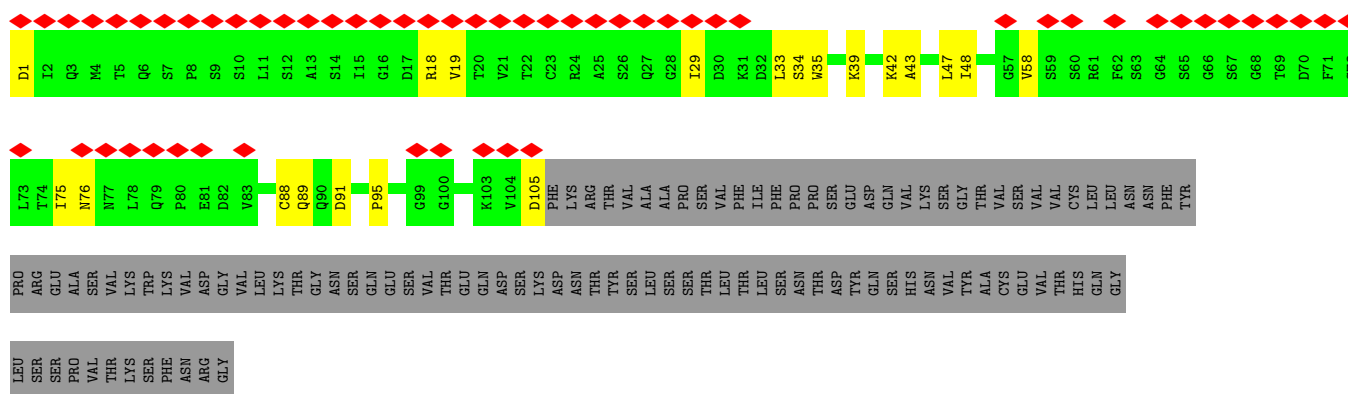
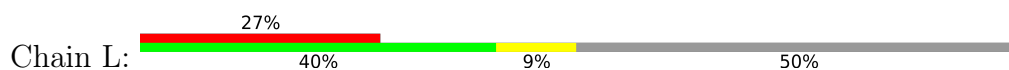




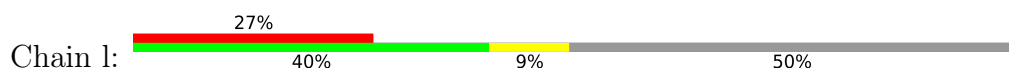
• Molecule 3: DFPH-a.15 Light chain



• Molecule 3: DFPH-a.15 Light chain

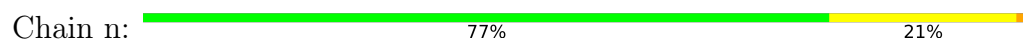


• Molecule 3: DFPH-a.15 Light chain

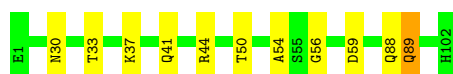
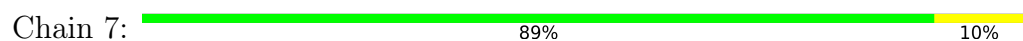




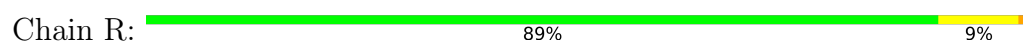
- Molecule 5: PGT122 Light chain



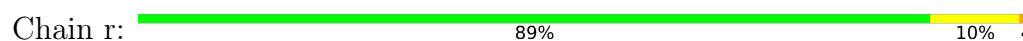
- Molecule 6: VRC03 Light chain



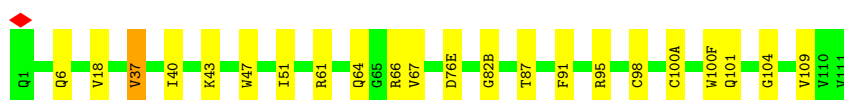
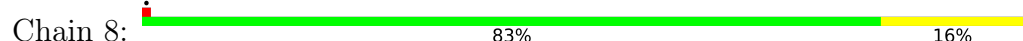
- Molecule 6: VRC03 Light chain



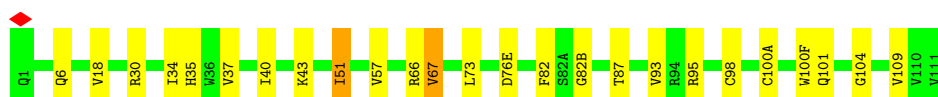
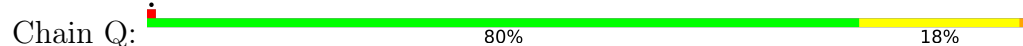
- Molecule 6: VRC03 Light chain



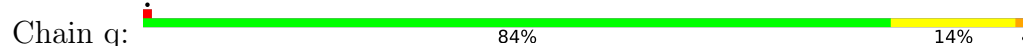
- Molecule 7: VRC03 Heavy chain



- Molecule 7: VRC03 Heavy chain

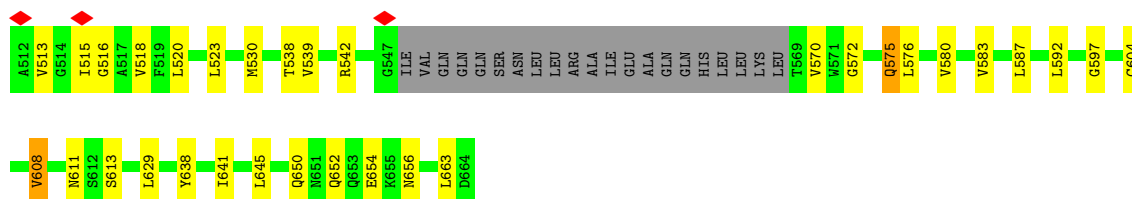


- Molecule 7: VRC03 Heavy chain

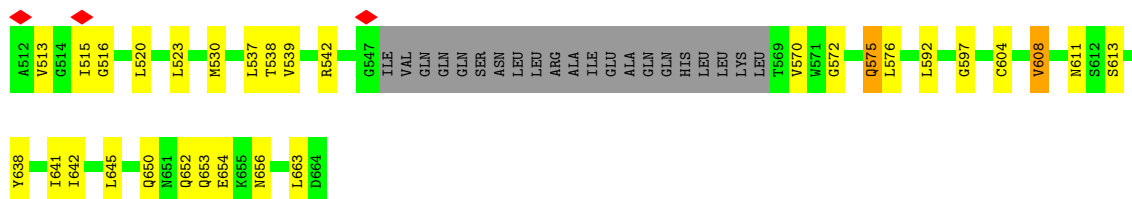




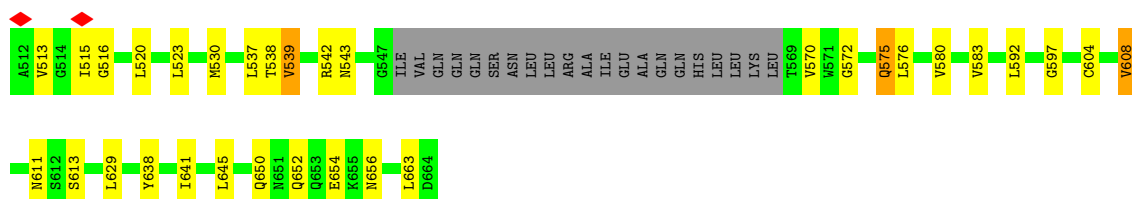
• Molecule 8: Envelope glycoprotein gp41



• Molecule 8: Envelope glycoprotein gp41



• Molecule 8: Envelope glycoprotein gp41



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



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





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

Chain G:  50% 50%



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

Chain I:  100%



- 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 O:  50% 50%



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

Chain U:  100%



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

Chain W:  50% 50%



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

Chain Y:  100%



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

Chain Z:  50% 50%



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

Chain a:  100%



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

Chain e:  100%

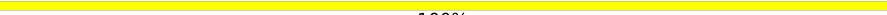


- Molecule 9: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain f:  50% 50%



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

Chain k:  100%



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

Chain p:  50% 50%



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

Chain t:  100%



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

Chain u:  50%  50%



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

Chain v:  100%



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

Chain x:  100%



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

Chain y:  100%



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

Chain 9:  100%



- Molecule 10: 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



- Molecule 10: 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



- Molecule 10: 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



- Molecule 10: 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



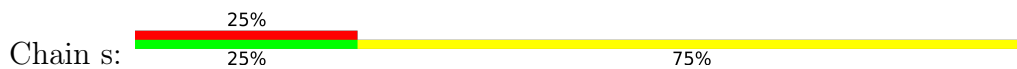
- Molecule 10: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 10: 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



- Molecule 10: 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



- Molecule 10: 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



- Molecule 10: 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



- 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



- 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



- 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



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

Chain P:  67% 33%



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

Chain g:  33% 67%

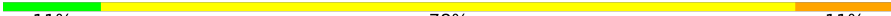


- Molecule 12: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain z:  67% 33%

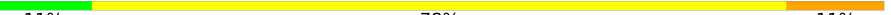


- Molecule 13: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:  11% 78% 11%



- Molecule 13: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain j:  11% 78% 11%



- Molecule 13: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain 1:  11% 89%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	56291	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 ($e^-/\text{\AA}^2$)	69.98	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	22500	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	3.572	Depositor
Minimum map value	-1.576	Depositor
Average map value	0.010	Depositor
Map value standard deviation	0.105	Depositor
Recommended contour level	0.688	Depositor
Map size (Å)	410.88, 410.88, 410.88	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.07, 1.07, 1.07	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	2	0.26	0/3638	0.51	0/4939
1	C	0.26	0/3638	0.50	0/4939
1	c	0.27	0/3638	0.52	0/4939
2	3	0.21	0/1003	0.48	0/1363
2	H	0.20	0/1003	0.48	0/1363
2	h	0.21	0/1003	0.48	0/1363
3	4	0.17	0/802	0.45	0/1088
3	L	0.17	0/802	0.45	0/1088
3	l	0.17	0/802	0.45	0/1088
4	5	0.19	0/1076	0.46	0/1465
4	M	0.20	0/1076	0.48	0/1465
4	m	0.20	0/1076	0.47	0/1465
5	6	0.27	0/826	0.49	0/1130
5	N	0.26	0/826	0.49	0/1130
5	n	0.26	0/826	0.49	0/1130
6	7	0.24	0/820	0.53	0/1107
6	R	0.25	0/820	0.52	0/1107
6	r	0.24	0/820	0.52	0/1107
7	8	0.26	0/1056	0.43	0/1439
7	Q	0.26	0/1056	0.44	0/1439
7	q	0.26	0/1056	0.43	0/1439
8	A	0.27	0/1052	0.52	0/1427
8	D	0.27	0/1052	0.53	0/1427
8	d	0.27	0/1052	0.52	0/1427
All	All	0.24	0/30819	0.49	0/41874

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
3	4	0	1
3	L	0	1
3	l	0	1
6	7	0	1
6	R	0	1
6	r	0	1
All	All	0	6

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	4	29	ILE	Peptide
6	7	89	GLN	Peptide
3	L	29	ILE	Peptide
6	R	89	GLN	Peptide
3	l	29	ILE	Peptide
6	r	89	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	3495	45	0
1	C	3564	0	3495	48	0
1	c	3564	0	3495	44	0
2	3	980	0	955	19	0
2	H	980	0	955	18	0
2	h	980	0	955	18	0
3	4	787	0	760	12	0
3	L	787	0	760	12	0
3	l	787	0	760	13	0
4	5	1047	0	1026	15	0
4	M	1047	0	1026	14	0
4	m	1047	0	1026	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	6	805	0	761	9	0
5	N	805	0	761	6	0
5	n	805	0	761	11	0
6	7	802	0	780	6	0
6	R	802	0	780	6	0
6	r	802	0	780	5	0
7	8	1023	0	971	12	0
7	Q	1023	0	971	14	0
7	q	1023	0	971	11	0
8	A	1034	0	1015	12	0
8	D	1034	0	1015	11	0
8	d	1034	0	1015	12	0
9	9	28	0	25	0	0
9	B	28	0	25	0	0
9	F	28	0	25	0	0
9	G	28	0	25	0	0
9	I	28	0	25	0	0
9	K	28	0	25	0	0
9	O	28	0	25	0	0
9	U	28	0	25	0	0
9	W	28	0	25	0	0
9	Y	28	0	25	0	0
9	Z	28	0	25	0	0
9	a	28	0	25	0	0
9	e	28	0	25	0	0
9	f	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	u	28	0	25	0	0
9	v	28	0	25	0	0
9	x	28	0	25	0	0
9	y	28	0	25	0	0
10	0	50	0	43	0	0
10	AA	50	0	43	1	0
10	E	50	0	43	0	0
10	S	50	0	43	0	0
10	V	50	0	43	1	0
10	X	50	0	43	0	0
10	i	50	0	43	0	0
10	o	50	0	43	1	0
10	s	50	0	43	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
11	J	61	0	52	0	0
11	b	61	0	52	0	0
11	w	61	0	52	0	0
12	P	39	0	34	0	0
12	g	39	0	34	0	0
12	z	39	0	34	0	0
13	1	105	0	88	0	0
13	T	105	0	88	1	0
13	j	105	0	88	2	0
14	2	42	0	39	1	0
14	C	42	0	39	0	0
14	c	42	0	39	0	0
All	All	31905	0	30840	364	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 (364) 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)
3:l:18:ARG:HA	3:l:75:ILE:O	1.73	0.86
3:L:18:ARG:HA	3:L:75:ILE:O	1.78	0.84
3:4:18:ARG:HA	3:4:75:ILE:O	1.77	0.83
1:c:39:TYR:O	1:c:494:LEU:HA	1.86	0.76
1:C:98:ASN:HD21	1:C:100:MET:HE2	1.57	0.69
4:m:52:HIS:HB3	4:m:56:ASP:HB3	1.76	0.68
4:5:52:HIS:HB3	4:5:56:ASP:HB3	1.76	0.68
1:2:39:TYR:O	1:2:494:LEU:HA	1.94	0.67
4:M:52:HIS:HB3	4:M:56:ASP:HB3	1.76	0.67
2:h:22:CYS:HB3	2:h:78:PHE:HB3	1.78	0.65
1:C:195:ASN:ND2	1:C:201:CYS:SG	2.69	0.65
1:c:195:ASN:ND2	1:c:201:CYS:SG	2.70	0.65
2:3:22:CYS:HB3	2:3:78:PHE:HB3	1.79	0.64
2:H:22:CYS:HB3	2:H:78:PHE:HB3	1.78	0.64
1:2:98:ASN:HD21	1:2:100:MET:HE2	1.62	0.64
1:2:195:ASN:ND2	1:2:201:CYS:SG	2.70	0.64
1:c:98:ASN:HD21	1:c:100:MET:HE2	1.62	0.64
1:C:95:MET:HE1	1:C:273:ARG:HD3	1.80	0.64
2:H:68:ILE:HB	2:H:81:ARG:HB2	1.81	0.63
2:h:68:ILE:HB	2:h:81:ARG:HB2	1.81	0.62
2:3:68:ILE:HB	2:3:81:ARG:HB2	1.80	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:95:MET:HE1	1:2:273:ARG:HD3	1.83	0.61
1:C:39:TYR:O	1:C:494:LEU:HA	2.01	0.61
7:Q:40:ILE:HB	7:Q:43:LYS:HB2	1.83	0.60
1:c:95:MET:HE1	1:c:273:ARG:HD3	1.82	0.60
1:2:104:MET:HE3	1:2:479:TRP:HB3	1.83	0.60
5:6:22:THR:HG22	5:6:72:THR:HG22	1.84	0.59
5:n:22:THR:HG22	5:n:72:THR:HG22	1.84	0.59
1:c:365:SER:OG	1:c:469:ARG:NH1	2.35	0.59
1:C:265:LEU:HD23	1:C:450:THR:HG23	1.85	0.59
5:N:22:THR:HG22	5:N:72:THR:HG22	1.85	0.59
1:c:195:ASN:OD1	1:c:195:ASN:N	2.32	0.59
3:l:39:LYS:HB2	3:l:42:LYS:HB2	1.85	0.59
7:8:40:ILE:HB	7:8:43:LYS:HB2	1.86	0.58
1:c:104:MET:HE3	1:c:479:TRP:HB3	1.85	0.58
3:l:91:ASP:OD1	3:l:91:ASP:N	2.37	0.58
3:L:39:LYS:HB2	3:L:42:LYS:HB2	1.84	0.58
3:L:34:SER:O	3:L:88:CYS:HA	2.05	0.57
1:c:499:THR:OG1	1:c:500:ARG:N	2.38	0.57
4:5:60:ASN:ND2	4:5:62:SER:OG	2.37	0.57
7:q:40:ILE:HB	7:q:43:LYS:HB2	1.86	0.56
2:h:21:THR:HA	2:h:78:PHE:O	2.04	0.56
6:r:44:ARG:NH2	6:r:56:GLY:O	2.38	0.56
1:C:499:THR:OG1	1:C:500:ARG:N	2.38	0.56
1:2:324:GLY:O	5:n:94:ARG:NH2	2.39	0.56
5:6:94:ARG:NH2	1:C:324:GLY:O	2.38	0.56
3:4:91:ASP:OD1	3:4:91:ASP:N	2.38	0.56
1:2:365:SER:OG	1:2:469:ARG:NH1	2.39	0.56
6:R:30:ASN:HB2	6:R:89:GLN:HE21	1.71	0.56
7:Q:87:THR:HA	7:Q:109:VAL:O	2.06	0.56
3:4:39:LYS:HB2	3:4:42:LYS:HB2	1.86	0.55
6:7:44:ARG:NH2	6:7:56:GLY:O	2.39	0.55
2:3:97:LYS:HA	2:3:100(H):ARG:HA	1.88	0.55
2:H:21:THR:HA	2:H:78:PHE:O	2.06	0.55
6:r:30:ASN:HB2	6:r:89:GLN:HE21	1.70	0.55
3:L:91:ASP:OD1	3:L:91:ASP:N	2.39	0.55
1:c:476:ARG:HA	1:c:479:TRP:HD1	1.72	0.55
4:m:100:ARG:NH1	4:m:100(A):ILE:O	2.39	0.55
7:8:66:ARG:NH2	7:8:82(B):GLY:O	2.39	0.55
1:2:265:LEU:HD23	1:2:450:THR:HG23	1.88	0.55
1:c:503:ARG:HH12	8:d:597:GLY:HA3	1.72	0.55
6:R:44:ARG:HH21	6:R:57:VAL:HG22	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:195:ASN:OD1	1:2:195:ASN:N	2.34	0.54
1:C:476:ARG:HA	1:C:479:TRP:HD1	1.72	0.54
2:h:97:LYS:HA	2:h:100(H):ARG:HA	1.89	0.54
7:q:87:THR:HA	7:q:109:VAL:O	2.07	0.54
8:A:608:VAL:HG11	8:A:645:LEU:HB3	1.89	0.54
1:C:270:VAL:HG11	1:C:345:VAL:HG12	1.89	0.54
4:5:38:ARG:HB3	4:5:48:ILE:HD11	1.89	0.54
4:m:30:ARG:HH21	4:m:73:LYS:HD3	1.72	0.54
8:D:608:VAL:HG11	8:D:645:LEU:HB3	1.90	0.54
2:3:21:THR:HA	2:3:78:PHE:O	2.07	0.54
3:l:33:LEU:HD11	3:l:88:CYS:HB2	1.90	0.54
1:2:476:ARG:HA	1:2:479:TRP:HD1	1.72	0.54
1:2:499:THR:OG1	1:2:500:ARG:N	2.40	0.54
7:8:76(E):ASP:OD1	7:8:76(E):ASP:N	2.32	0.54
1:C:116:LEU:HB3	1:C:202:THR:HG21	1.89	0.54
3:4:34:SER:O	3:4:88:CYS:HA	2.08	0.54
4:5:72:ASP:O	4:5:76:ASN:N	2.40	0.54
4:M:30:ARG:HH21	4:M:73:LYS:HD3	1.73	0.54
1:c:270:VAL:HG11	1:c:345:VAL:HG12	1.90	0.54
1:2:67:ASN:OD1	1:2:67:ASN:N	2.41	0.53
1:2:270:VAL:HG11	1:2:345:VAL:HG12	1.89	0.53
1:c:265:LEU:HD23	1:c:450:THR:HG23	1.89	0.53
3:4:33:LEU:HD11	3:4:88:CYS:HB2	1.90	0.53
2:H:97:LYS:HA	2:H:100(H):ARG:HA	1.89	0.53
2:h:87:THR:HG23	2:h:110:THR:HA	1.90	0.53
4:m:72:ASP:O	4:m:76:ASN:N	2.41	0.53
3:L:33:LEU:HD11	3:L:88:CYS:HB2	1.91	0.53
4:m:60:ASN:ND2	4:m:62:SER:OG	2.41	0.53
4:M:72:ASP:O	4:M:76:ASN:N	2.42	0.53
1:2:38:VAL:HG23	8:A:604:CYS:HB3	1.91	0.53
6:r:37:LYS:NZ	6:r:41:GLN:O	2.42	0.53
1:C:256:SER:O	1:C:478:ASN:ND2	2.38	0.53
2:h:19:SER:HA	2:h:80:LEU:O	2.08	0.53
1:2:33:ASN:OD1	1:2:33:ASN:N	2.40	0.52
1:2:371:VAL:HG23	1:2:473:GLY:H	1.74	0.52
7:8:64:GLN:NE2	1:C:457:ASP:OD2	2.41	0.52
1:C:503:ARG:HH12	8:D:597:GLY:HA3	1.74	0.52
4:m:38:ARG:HB3	4:m:48:ILE:HD11	1.91	0.52
1:C:97:LYS:NZ	10:o:2:NAG:O7	2.42	0.52
1:c:116:LEU:HB3	1:c:202:THR:HG21	1.90	0.52
1:C:365:SER:OG	1:C:469:ARG:NH1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:97:LYS:NZ	10:V:2:NAG:O7	2.42	0.52
1:C:104:MET:HE3	1:C:479:TRP:HB3	1.92	0.52
3:l:34:SER:O	3:l:88:CYS:HA	2.09	0.52
2:H:84:ALA:HA	2:H:111:VAL:HB	1.92	0.52
4:5:30:ARG:HH21	4:5:73:LYS:HD3	1.74	0.52
2:h:84:ALA:HA	2:h:111:VAL:HB	1.91	0.52
2:H:87:THR:HG23	2:H:110:THR:HA	1.92	0.51
4:M:60:ASN:ND2	4:M:62:SER:OG	2.44	0.51
1:c:216:HIS:ND1	1:c:248:THR:O	2.35	0.51
8:d:608:VAL:HG11	8:d:645:LEU:HB3	1.91	0.51
2:3:19:SER:HA	2:3:80:LEU:O	2.10	0.51
1:C:38:VAL:HG23	8:D:604:CYS:HB3	1.92	0.51
2:3:24:VAL:O	2:3:76:ASN:ND2	2.43	0.51
8:d:572:GLY:O	8:d:576:LEU:HB2	2.10	0.51
8:d:650:GLN:HE21	8:d:654:GLU:HG2	1.75	0.51
7:Q:66:ARG:NH2	7:Q:82(B):GLY:O	2.44	0.51
1:c:174:SER:OG	1:c:175:LEU:N	2.43	0.51
2:3:33:TYR:HB3	2:3:95:ARG:HB3	1.92	0.51
2:H:19:SER:HA	2:H:80:LEU:O	2.10	0.51
2:h:33:TYR:HB3	2:h:95:ARG:HB3	1.91	0.51
2:3:84:ALA:HA	2:3:111:VAL:HB	1.92	0.51
5:6:38:GLN:NE2	5:6:39:ARG:O	2.44	0.51
1:c:371:VAL:HG23	1:c:473:GLY:H	1.75	0.51
1:2:503:ARG:HH12	8:A:597:GLY:HA3	1.76	0.50
6:R:44:ARG:NH2	6:R:56:GLY:O	2.44	0.50
2:3:20:LEU:O	2:3:79:SER:HA	2.12	0.50
1:C:113:ASP:OD2	1:C:429:ARG:NH2	2.45	0.50
1:c:33:ASN:OD1	1:c:33:ASN:N	2.45	0.50
7:8:6:GLN:HE21	7:8:104:GLY:HA3	1.75	0.50
1:c:38:VAL:HG23	8:d:604:CYS:HB3	1.93	0.50
7:q:95:ARG:HA	7:q:100(F):TRP:HA	1.92	0.50
2:h:20:LEU:O	2:h:79:SER:HA	2.12	0.50
7:q:66:ARG:NH2	7:q:82(B):GLY:O	2.44	0.50
1:C:174:SER:OG	1:C:175:LEU:N	2.44	0.50
1:c:298:ARG:NH1	1:c:381:GLU:OE2	2.44	0.50
2:3:87:THR:HG23	2:3:110:THR:HA	1.94	0.50
2:h:24:VAL:O	2:h:76:ASN:ND2	2.44	0.50
4:M:38:ARG:HB3	4:M:48:ILE:HD11	1.94	0.50
5:n:38:GLN:NE2	5:n:39:ARG:O	2.45	0.49
6:7:37:LYS:NZ	6:7:41:GLN:O	2.45	0.49
2:H:24:VAL:O	2:H:76:ASN:ND2	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:q:6:GLN:HE21	7:q:104:GLY:HA3	1.76	0.49
2:H:33:TYR:HB3	2:H:95:ARG:HB3	1.93	0.49
1:c:97:LYS:NZ	10:AA:2:NAG:O7	2.45	0.49
1:c:304:ARG:N	1:c:440:GLN:OE1	2.45	0.49
1:2:459:GLY:HA3	7:q:61:ARG:HH11	1.78	0.49
7:8:87:THR:HA	7:8:109:VAL:O	2.13	0.49
4:M:28:LEU:HA	4:M:76:ASN:HD21	1.77	0.49
1:C:33:ASN:OD1	1:C:33:ASN:N	2.41	0.49
6:R:33:THR:HB	6:R:88:GLN:HB3	1.94	0.49
2:H:20:LEU:O	2:H:79:SER:HA	2.13	0.48
7:8:98:CYS:SG	7:8:100(A):CYS:N	2.86	0.48
1:C:299:PRO:HG2	1:C:327:ARG:HB2	1.95	0.48
8:D:516:GLY:HA3	2:H:96:ALA:HA	1.96	0.48
1:c:188:ASN:OD1	1:c:188:ASN:N	2.47	0.48
7:Q:76(E):ASP:OD1	7:Q:76(E):ASP:N	2.34	0.48
8:A:650:GLN:HE21	8:A:654:GLU:HG2	1.78	0.48
5:6:51:ASN:ND2	13:j:5:MAN:O4	2.45	0.47
5:N:35:TRP:HB2	5:N:48:ILE:HB	1.95	0.47
5:n:50:ASN:ND2	13:T:5:MAN:O6	2.45	0.47
1:2:304:ARG:N	1:2:440:GLN:OE1	2.47	0.47
4:5:28:LEU:HA	4:5:76:ASN:HD21	1.79	0.47
1:C:67:ASN:OD1	1:C:67:ASN:N	2.42	0.47
1:2:65:LYS:HE2	1:2:65:LYS:HB3	1.70	0.47
3:4:1:ASP:HA	3:4:95:PRO:HD2	1.97	0.47
8:d:530:MET:HE3	8:d:530:MET:HB2	1.74	0.47
6:7:30:ASN:HB2	6:7:89:GLN:HE21	1.80	0.47
1:C:277:ILE:HG13	1:C:352:HIS:HD2	1.80	0.47
1:c:272:ILE:HG22	1:c:286:VAL:HG22	1.95	0.47
3:4:19:VAL:HG22	3:4:75:ILE:HB	1.97	0.47
8:D:650:GLN:HE21	8:D:654:GLU:HG2	1.80	0.47
7:Q:67:VAL:HG23	7:Q:82:PHE:HD1	1.79	0.47
3:l:35:TRP:HB2	3:l:48:ILE:HB	1.95	0.47
4:m:52:HIS:ND1	4:m:53:ASP:O	2.48	0.47
1:C:270:VAL:HB	1:C:348:GLN:HG3	1.96	0.47
1:C:371:VAL:HG23	1:C:473:GLY:H	1.80	0.47
1:C:377:ASN:OD1	1:C:381:GLU:N	2.48	0.47
3:4:35:TRP:HB2	3:4:48:ILE:HB	1.96	0.47
3:l:1:ASP:HA	3:l:95:PRO:HD2	1.97	0.47
1:2:277:ILE:HG13	1:2:352:HIS:HD2	1.79	0.46
4:5:100(Q):ASP:OD1	4:5:100(Q):ASP:N	2.47	0.46
2:3:96:ALA:HA	8:A:516:GLY:HA3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:65:LYS:HB3	1:C:65:LYS:HE2	1.71	0.46
4:m:7:SER:OG	4:m:21:THR:OG1	2.33	0.46
3:L:19:VAL:HG22	3:L:75:ILE:HB	1.97	0.46
3:L:35:TRP:HB2	3:L:48:ILE:HB	1.96	0.46
1:2:216:HIS:ND1	1:2:248:THR:O	2.39	0.46
6:r:59:ASP:OD1	6:r:59:ASP:N	2.48	0.46
2:h:38:ARG:HB3	2:h:48:ILE:HD11	1.97	0.46
4:m:88:ALA:HB3	4:m:90:TYR:HE1	1.80	0.46
5:n:85:ASP:N	5:n:85:ASP:OD1	2.49	0.46
4:5:88:ALA:HB3	4:5:90:TYR:HE1	1.81	0.46
7:8:47:TRP:HZ3	1:C:458:GLY:HA3	1.81	0.46
1:C:272:ILE:HG22	1:C:286:VAL:HG22	1.97	0.46
8:D:572:GLY:O	8:D:576:LEU:HB2	2.16	0.46
7:Q:6:GLN:HE21	7:Q:104:GLY:HA3	1.80	0.46
1:c:256:SER:O	1:c:478:ASN:ND2	2.44	0.46
1:2:272:ILE:HG22	1:2:286:VAL:HG22	1.97	0.45
2:3:49:GLY:HA2	2:3:59:HIS:HA	1.97	0.45
8:d:516:GLY:HA3	2:h:96:ALA:HA	1.98	0.45
2:h:49:GLY:HA2	2:h:59:HIS:HA	1.98	0.45
7:8:37:VAL:HG12	7:8:91:PHE:HB2	1.98	0.45
1:c:457:ASP:HB2	1:c:467:THR:HG23	1.99	0.45
2:3:38:ARG:HB3	2:3:48:ILE:HD11	1.99	0.45
1:C:50:THR:OG1	1:C:51:THR:N	2.49	0.45
1:c:67:ASN:OD1	1:c:67:ASN:N	2.46	0.45
1:c:288:PHE:HZ	1:c:449:ILE:HG22	1.80	0.45
7:q:98:CYS:SG	7:q:100(A):CYS:N	2.90	0.45
1:2:116:LEU:HD21	1:2:434:MET:HE2	1.98	0.45
1:2:174:SER:OG	1:2:175:LEU:N	2.49	0.45
4:5:22:CYS:HB3	4:5:78:VAL:HB	1.99	0.45
3:L:105:ASP:N	3:L:105:ASP:OD1	2.50	0.45
4:M:88:ALA:HB3	4:M:90:TYR:HE1	1.81	0.45
7:q:51:ILE:HB	7:q:57:VAL:HG12	1.98	0.45
5:6:50:ASN:ND2	13:j:5:MAN:O6	2.44	0.45
1:C:195:ASN:OD1	1:C:195:ASN:N	2.37	0.45
1:C:257:THR:OG1	1:C:258:GLN:N	2.49	0.45
1:C:476:ARG:HA	1:C:479:TRP:CD1	2.51	0.45
5:6:33:VAL:HG12	5:6:51:ASN:HA	1.99	0.45
8:A:652:GLN:OE1	8:A:656:ASN:ND2	2.50	0.45
1:2:257:THR:OG1	1:2:258:GLN:N	2.48	0.45
2:H:49:GLY:HA2	2:H:59:HIS:HA	1.98	0.45
8:D:611:ASN:HD21	8:D:613:SER:HB3	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:M:52:HIS:ND1	4:M:53:ASP:O	2.50	0.45
7:Q:51:ILE:HB	7:Q:57:VAL:HG12	1.99	0.45
1:2:113:ASP:OD2	1:2:429:ARG:NH2	2.51	0.44
1:C:132:THR:OG1	1:C:133:ASN:N	2.50	0.44
3:l:19:VAL:HG22	3:l:75:ILE:HB	1.99	0.44
4:5:100:ARG:NH1	4:5:100(A):ILE:O	2.50	0.44
1:C:92:GLU:HA	1:C:238:PRO:HA	1.99	0.44
3:4:105:ASP:N	3:4:105:ASP:OD1	2.50	0.44
1:2:116:LEU:HB3	1:2:202:THR:HG21	1.99	0.44
1:c:299:PRO:HG2	1:c:327:ARG:HB2	1.99	0.44
8:d:652:GLN:OE1	8:d:656:ASN:ND2	2.51	0.44
1:2:132:THR:OG1	1:2:133:ASN:N	2.51	0.44
1:2:299:PRO:HG2	1:2:327:ARG:HB2	1.98	0.44
5:6:85:ASP:OD1	5:6:85:ASP:N	2.51	0.44
6:7:59:ASP:OD1	6:7:59:ASP:N	2.50	0.44
1:2:166:ARG:HB2	1:c:127:VAL:HG23	2.00	0.44
1:2:50:THR:OG1	1:2:51:THR:N	2.51	0.44
1:C:358:ILE:HB	1:C:465:THR:HG22	2.00	0.44
8:A:575:GLN:HE21	8:A:575:GLN:HB3	1.51	0.43
4:M:100(Q):ASP:OD1	4:M:100(Q):ASP:N	2.43	0.43
1:2:256:SER:O	1:2:478:ASN:ND2	2.39	0.43
8:A:580:VAL:HA	8:A:583:VAL:HG22	2.00	0.43
8:D:575:GLN:HE21	8:D:575:GLN:HB3	1.54	0.43
3:l:105:ASP:N	3:l:105:ASP:OD1	2.50	0.43
8:A:572:GLY:O	8:A:576:LEU:HB2	2.18	0.43
5:N:31:ARG:HG2	5:N:92:ASP:HA	2.00	0.43
1:c:166:ARG:HA	1:c:166:ARG:HD3	1.80	0.43
3:4:47:LEU:HD23	3:4:58:VAL:HG11	2.01	0.43
7:Q:93:VAL:HB	7:Q:100(F):TRP:HB3	1.99	0.43
5:N:85:ASP:OD1	5:N:85:ASP:N	2.52	0.43
2:h:60:ASN:ND2	2:h:62:SER:OG	2.52	0.43
5:N:46:LEU:HD21	5:N:49:TYR:HB3	1.99	0.43
4:5:7:SER:OG	4:5:21:THR:OG1	2.36	0.43
4:5:52:HIS:ND1	4:5:53:ASP:O	2.51	0.43
1:c:270:VAL:HB	1:c:348:GLN:HG3	2.01	0.43
4:m:100(Q):ASP:N	4:m:100(Q):ASP:OD1	2.43	0.43
5:n:35:TRP:HB2	5:n:48:ILE:HB	2.01	0.43
1:2:430:ILE:H	1:2:430:ILE:HG13	1.63	0.43
3:L:34:SER:HB2	3:L:89:GLN:HB3	2.01	0.43
4:M:22:CYS:HB3	4:M:78:VAL:HB	2.00	0.43
1:c:65:LYS:HB3	1:c:65:LYS:HE2	1.78	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:m:28:LEU:HA	4:m:76:ASN:HD21	1.83	0.43
1:2:92:GLU:HA	1:2:238:PRO:HA	2.01	0.42
8:d:539:VAL:O	8:d:543:ASN:ND2	2.51	0.42
1:2:354:GLY:HA2	14:2:623:NAG:H82	2.00	0.42
2:H:47:TRP:NE1	2:H:49:GLY:O	2.52	0.42
1:c:277:ILE:HG13	1:c:352:HIS:HD2	1.85	0.42
5:n:58:ILE:HD12	5:n:59:PRO:HD2	2.01	0.42
2:3:20:LEU:HD11	2:3:109:VAL:HG21	2.00	0.42
2:H:20:LEU:HD11	2:H:109:VAL:HG21	2.00	0.42
2:h:20:LEU:HD11	2:h:109:VAL:HG21	2.01	0.42
1:2:188:ASN:OD1	1:2:188:ASN:N	2.46	0.42
7:Q:95:ARG:HA	7:Q:100(F):TRP:HA	2.01	0.42
5:n:46:LEU:HD21	5:n:49:TYR:HB3	2.01	0.42
5:6:94:ARG:O	1:C:137:ASN:N	2.53	0.42
7:8:61:ARG:HE	1:C:459:GLY:HA2	1.85	0.42
1:C:475:MET:HE3	1:C:475:MET:HB3	1.96	0.42
5:N:58:ILE:HD12	5:N:59:PRO:HD2	2.01	0.42
7:Q:98:CYS:SG	7:Q:100(A):CYS:N	2.92	0.42
1:c:360:ARG:HG3	1:c:467:THR:HB	2.01	0.42
4:m:72:ASP:O	4:m:76:ASN:CA	2.67	0.42
7:q:37:VAL:HG12	7:q:91:PHE:HB2	2.02	0.42
4:5:7:SER:HG	4:5:21:THR:HG1	1.63	0.42
1:c:92:GLU:HA	1:c:238:PRO:HA	2.00	0.42
8:A:530:MET:HE3	8:A:530:MET:HB2	1.75	0.42
2:H:105:PRO:HA	3:L:43:ALA:HB2	2.02	0.42
1:c:350:ARG:NH2	1:c:396:ILE:O	2.46	0.42
7:q:76(E):ASP:OD1	7:q:76(E):ASP:N	2.38	0.42
7:8:95:ARG:HA	7:8:100(F):TRP:HA	2.01	0.42
3:L:1:ASP:HA	3:L:95:PRO:HD2	2.00	0.42
7:Q:35:HIS:HB2	7:Q:93:VAL:HG23	2.01	0.42
1:2:346:VAL:HG21	1:2:395:TRP:CG	2.54	0.42
1:C:121:LYS:HB2	1:C:121:LYS:HE2	1.80	0.42
7:Q:101:GLN:HE22	6:R:54:ALA:HB1	1.84	0.42
2:3:47:TRP:NE1	2:3:49:GLY:O	2.53	0.42
1:C:346:VAL:HG21	1:C:395:TRP:CG	2.55	0.41
4:M:72:ASP:O	4:M:76:ASN:CA	2.68	0.41
8:d:580:VAL:HA	8:d:583:VAL:HG22	2.01	0.41
8:d:611:ASN:HD21	8:d:613:SER:HB3	1.85	0.41
2:3:99:TYR:HE2	2:3:100(B):ALA:HA	1.85	0.41
5:6:58:ILE:HD12	5:6:59:PRO:HD2	2.01	0.41
8:A:611:ASN:HD21	8:A:613:SER:HB3	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:496:VAL:HG21	8:D:642:ILE:HG21	2.02	0.41
3:L:47:LEU:HD23	3:L:58:VAL:HG11	2.02	0.41
3:l:33:LEU:HB3	3:l:51:ALA:HB2	2.02	0.41
7:Q:30:ARG:HE	7:Q:73:LEU:HD13	1.84	0.41
6:R:59:ASP:OD1	6:R:59:ASP:N	2.48	0.41
1:c:121:LYS:HE2	1:c:121:LYS:HB2	1.82	0.41
4:m:36:TRP:HB3	4:m:48:ILE:HD12	2.03	0.41
2:h:99:TYR:HE2	2:h:100(B):ALA:HA	1.85	0.41
2:h:105:PRO:HA	3:l:43:ALA:HB2	2.03	0.41
7:q:67:VAL:HG23	7:q:82:PHE:HD1	1.85	0.41
1:2:457:ASP:HB2	1:2:467:THR:HG23	2.02	0.41
2:3:105:PRO:HA	3:4:43:ALA:HB2	2.02	0.41
1:C:338:TRP:CZ2	1:C:390:LEU:HB3	2.55	0.41
3:l:47:LEU:HD23	3:l:58:VAL:HG11	2.01	0.41
6:7:54:ALA:HB1	7:8:101:GLN:HE22	1.86	0.41
8:D:652:GLN:OE1	8:D:656:ASN:ND2	2.54	0.41
4:M:39:GLN:NE2	4:M:43:LYS:O	2.38	0.41
1:C:266:ALA:HB2	1:C:287:GLN:HG2	2.01	0.41
1:c:113:ASP:OD2	1:c:429:ARG:NH2	2.54	0.41
4:m:36:TRP:HA	4:m:91:TYR:O	2.21	0.41
5:n:33:VAL:HG12	5:n:51:ASN:HA	2.03	0.41
1:2:297:THR:HA	1:2:443:ILE:O	2.21	0.41
4:5:72:ASP:O	4:5:76:ASN:CA	2.68	0.41
2:H:38:ARG:HB3	2:H:48:ILE:HD11	2.03	0.41
1:c:398:ASN:OD1	1:c:398:ASN:N	2.52	0.41
3:l:34:SER:HB2	3:l:89:GLN:HB3	2.02	0.41
5:n:31:ARG:HG2	5:n:92:ASP:HA	2.03	0.41
1:2:227:LYS:HE2	1:2:227:LYS:HB3	1.93	0.41
2:H:99:TYR:HE2	2:H:100(B):ALA:HA	1.86	0.41
1:c:385:CYS:HA	1:c:418:CYS:HA	2.03	0.41
8:d:575:GLN:HE21	8:d:575:GLN:HB3	1.53	0.41
2:h:50:GLY:H	2:h:69:ILE:HD13	1.86	0.41
1:C:166:ARG:HA	1:C:166:ARG:HD3	1.81	0.41
1:C:216:HIS:ND1	1:C:248:THR:O	2.48	0.41
2:H:82:LEU:HG	2:H:82(C):VAL:HG22	2.03	0.41
4:m:22:CYS:HB3	4:m:78:VAL:HB	2.03	0.41
1:2:288:PHE:HZ	1:2:449:ILE:HG22	1.86	0.40
2:3:82:LEU:HG	2:3:82(C):VAL:HG22	2.03	0.40
6:7:33:THR:HB	6:7:88:GLN:HB3	2.03	0.40
1:C:188:ASN:OD1	1:C:188:ASN:N	2.51	0.40
7:Q:34:ILE:HD13	7:Q:34:ILE:HA	1.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:m:34:TRP:HB2	4:m:51:VAL:HG13	2.03	0.40
5:n:19:ALA:HB3	5:n:75:ILE:HB	2.03	0.40
2:3:98:ILE:HA	8:A:518:VAL:HG22	2.02	0.40
3:4:34:SER:HB2	3:4:89:GLN:HB3	2.03	0.40
1:c:63:THR:O	1:c:67:ASN:ND2	2.45	0.40
1:c:266:ALA:HB2	1:c:287:GLN:HG2	2.03	0.40
1:2:476:ARG:HA	1:2:479:TRP:CD1	2.54	0.40
4:M:41:LEU:H	4:M:41:LEU:HG	1.70	0.40
6:r:33:THR:HB	6:r:88:GLN:HB3	2.04	0.40
1:2:398:ASN:OD1	1:2:398:ASN:N	2.53	0.40
4:M:34:TRP:HB2	4:M:51:VAL:HG13	2.02	0.40
1:c:168:LYS:HD3	1:c:168:LYS:HA	1.80	0.40
1:c:476:ARG:HA	1:c:479:TRP:CD1	2.55	0.40
4:5:29:VAL:HG21	4:5:71:LEU:HB3	2.03	0.40
1:C:503:ARG:NH2	8:D:653:GLN:OE1	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	2	447/473 (94%)	428 (96%)	19 (4%)	0	100	100
1	C	447/473 (94%)	426 (95%)	21 (5%)	0	100	100
1	c	447/473 (94%)	426 (95%)	21 (5%)	0	100	100
2	3	125/224 (56%)	119 (95%)	6 (5%)	0	100	100
2	H	125/224 (56%)	120 (96%)	5 (4%)	0	100	100
2	h	125/224 (56%)	119 (95%)	6 (5%)	0	100	100
3	4	103/212 (49%)	94 (91%)	9 (9%)	0	100	100
3	L	103/212 (49%)	94 (91%)	9 (9%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	l	103/212 (49%)	94 (91%)	9 (9%)	0	100	100
4	5	130/132 (98%)	123 (95%)	7 (5%)	0	100	100
4	M	130/132 (98%)	125 (96%)	5 (4%)	0	100	100
4	m	130/132 (98%)	124 (95%)	6 (5%)	0	100	100
5	6	103/105 (98%)	95 (92%)	8 (8%)	0	100	100
5	N	103/105 (98%)	94 (91%)	9 (9%)	0	100	100
5	n	103/105 (98%)	96 (93%)	7 (7%)	0	100	100
6	7	100/102 (98%)	94 (94%)	6 (6%)	0	100	100
6	R	100/102 (98%)	94 (94%)	6 (6%)	0	100	100
6	r	100/102 (98%)	94 (94%)	6 (6%)	0	100	100
7	8	126/128 (98%)	122 (97%)	4 (3%)	0	100	100
7	Q	126/128 (98%)	122 (97%)	4 (3%)	0	100	100
7	q	126/128 (98%)	122 (97%)	4 (3%)	0	100	100
8	A	128/153 (84%)	120 (94%)	8 (6%)	0	100	100
8	D	128/153 (84%)	121 (94%)	7 (6%)	0	100	100
8	d	128/153 (84%)	121 (94%)	7 (6%)	0	100	100
All	All	3786/4587 (82%)	3587 (95%)	199 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	2	405/422 (96%)	369 (91%)	36 (9%)	9	29
1	C	405/422 (96%)	371 (92%)	34 (8%)	10	31
1	c	405/422 (96%)	367 (91%)	38 (9%)	8	27
2	3	111/197 (56%)	105 (95%)	6 (5%)	20	43
2	H	111/197 (56%)	105 (95%)	6 (5%)	20	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	h	111/197 (56%)	105 (95%)	6 (5%)	20	43
3	4	90/188 (48%)	89 (99%)	1 (1%)	65	73
3	L	90/188 (48%)	89 (99%)	1 (1%)	65	73
3	l	90/188 (48%)	88 (98%)	2 (2%)	45	64
4	5	116/116 (100%)	108 (93%)	8 (7%)	14	37
4	M	116/116 (100%)	109 (94%)	7 (6%)	17	40
4	m	116/116 (100%)	108 (93%)	8 (7%)	14	37
5	6	88/88 (100%)	81 (92%)	7 (8%)	11	32
5	N	88/88 (100%)	81 (92%)	7 (8%)	11	32
5	n	88/88 (100%)	81 (92%)	7 (8%)	11	32
6	7	86/86 (100%)	85 (99%)	1 (1%)	63	73
6	R	86/86 (100%)	83 (96%)	3 (4%)	32	53
6	r	86/86 (100%)	84 (98%)	2 (2%)	44	64
7	8	108/108 (100%)	104 (96%)	4 (4%)	30	51
7	Q	108/108 (100%)	104 (96%)	4 (4%)	30	51
7	q	108/108 (100%)	104 (96%)	4 (4%)	30	51
8	A	110/129 (85%)	94 (86%)	16 (14%)	3	15
8	D	110/129 (85%)	94 (86%)	16 (14%)	3	15
8	d	110/129 (85%)	94 (86%)	16 (14%)	3	15
All	All	3342/4002 (84%)	3102 (93%)	240 (7%)	15	35

All (240) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	2	34	LEU
1	2	38	VAL
1	2	63	THR
1	2	67	ASN
1	2	74	CYS
1	2	75	VAL
1	2	101	VAL
1	2	111	LEU
1	2	125	LEU
1	2	138	ILE
1	2	162	THR

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Mol	Chain	Res	Type
1	2	182	VAL
1	2	195	ASN
1	2	198	THR
1	2	201	CYS
1	2	207	LYS
1	2	244	THR
1	2	254	VAL
1	2	260	LEU
1	2	294	ILE
1	2	303	THR
1	2	305	LYS
1	2	334	SER
1	2	357	THR
1	2	373	THR
1	2	374	HIS
1	2	375	SER
1	2	394	THR
1	2	415	THR
1	2	416	LEU
1	2	423	ILE
1	2	430	ILE
1	2	467	THR
1	2	488	VAL
1	2	496	VAL
1	2	499	THR
2	3	6	VAL
2	3	29	THR
2	3	35	SER
2	3	73	THR
2	3	95	ARG
2	3	98	ILE
3	4	76	ASN
4	5	41	LEU
4	5	57	THR
4	5	67	VAL
4	5	78	VAL
4	5	82	LEU
4	5	82(A)	THR
4	5	82(C)	VAL
4	5	109	VAL
5	6	12	SER
5	6	13	VAL

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Mol	Chain	Res	Type
5	6	34	ILE
5	6	58	ILE
5	6	85	ASP
5	6	105	ILE
5	6	106	VAL
6	7	50	THR
7	8	18	VAL
7	8	37	VAL
7	8	51	ILE
7	8	67	VAL
8	A	513	VAL
8	A	515	ILE
8	A	520	LEU
8	A	523	LEU
8	A	538	THR
8	A	539	VAL
8	A	542	ARG
8	A	570	VAL
8	A	575	GLN
8	A	587	LEU
8	A	592	LEU
8	A	608	VAL
8	A	629	LEU
8	A	638	TYR
8	A	641	ILE
8	A	663	LEU
1	C	34	LEU
1	C	38	VAL
1	C	63	THR
1	C	67	ASN
1	C	74	CYS
1	C	75	VAL
1	C	100	MET
1	C	101	VAL
1	C	111	LEU
1	C	122	LEU
1	C	125	LEU
1	C	138	ILE
1	C	195	ASN
1	C	198	THR
1	C	201	CYS
1	C	207	LYS

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Mol	Chain	Res	Type
1	C	244	THR
1	C	254	VAL
1	C	260	LEU
1	C	305	LYS
1	C	334	SER
1	C	357	THR
1	C	370	GLU
1	C	373	THR
1	C	374	HIS
1	C	375	SER
1	C	394	THR
1	C	396	ILE
1	C	416	LEU
1	C	430	ILE
1	C	467	THR
1	C	488	VAL
1	C	496	VAL
1	C	499	THR
8	D	513	VAL
8	D	515	ILE
8	D	520	LEU
8	D	523	LEU
8	D	530	MET
8	D	537	LEU
8	D	538	THR
8	D	539	VAL
8	D	542	ARG
8	D	570	VAL
8	D	575	GLN
8	D	592	LEU
8	D	608	VAL
8	D	638	TYR
8	D	641	ILE
8	D	663	LEU
2	H	6	VAL
2	H	29	THR
2	H	35	SER
2	H	73	THR
2	H	95	ARG
2	H	98	ILE
3	L	76	ASN
4	M	41	LEU

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Mol	Chain	Res	Type
4	M	57	THR
4	M	78	VAL
4	M	82	LEU
4	M	82(A)	THR
4	M	82(C)	VAL
4	M	109	VAL
5	N	12	SER
5	N	13	VAL
5	N	34	ILE
5	N	58	ILE
5	N	85	ASP
5	N	105	ILE
5	N	106	VAL
7	Q	18	VAL
7	Q	37	VAL
7	Q	51	ILE
7	Q	67	VAL
6	R	26	SER
6	R	44	ARG
6	R	50	THR
1	c	34	LEU
1	c	38	VAL
1	c	63	THR
1	c	67	ASN
1	c	74	CYS
1	c	75	VAL
1	c	100	MET
1	c	101	VAL
1	c	104	MET
1	c	111	LEU
1	c	125	LEU
1	c	138	ILE
1	c	195	ASN
1	c	198	THR
1	c	201	CYS
1	c	207	LYS
1	c	244	THR
1	c	254	VAL
1	c	260	LEU
1	c	294	ILE
1	c	297	THR
1	c	303	THR

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Mol	Chain	Res	Type
1	c	305	LYS
1	c	334	SER
1	c	346	VAL
1	c	357	THR
1	c	373	THR
1	c	374	HIS
1	c	375	SER
1	c	394	THR
1	c	416	LEU
1	c	430	ILE
1	c	445	CYS
1	c	448	ASN
1	c	467	THR
1	c	488	VAL
1	c	496	VAL
1	c	499	THR
8	d	513	VAL
8	d	515	ILE
8	d	520	LEU
8	d	523	LEU
8	d	537	LEU
8	d	538	THR
8	d	539	VAL
8	d	542	ARG
8	d	570	VAL
8	d	575	GLN
8	d	592	LEU
8	d	608	VAL
8	d	629	LEU
8	d	638	TYR
8	d	641	ILE
8	d	663	LEU
2	h	6	VAL
2	h	29	THR
2	h	35	SER
2	h	73	THR
2	h	95	ARG
2	h	98	ILE
3	l	76	ASN
3	l	91	ASP
4	m	41	LEU
4	m	57	THR

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Mol	Chain	Res	Type
4	m	67	VAL
4	m	78	VAL
4	m	82	LEU
4	m	82(A)	THR
4	m	82(C)	VAL
4	m	109	VAL
5	n	12	SER
5	n	13	VAL
5	n	34	ILE
5	n	58	ILE
5	n	85	ASP
5	n	105	ILE
5	n	106	VAL
7	q	18	VAL
7	q	37	VAL
7	q	51	ILE
7	q	67	VAL
6	r	50	THR
6	r	84	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
1	2	98	ASN
1	2	114	GLN
1	2	203	GLN
1	2	315	GLN
1	2	352	HIS
2	3	60	ASN
3	4	79	GLN
4	5	60	ASN
6	7	89	GLN
7	8	3	GLN
7	8	6	GLN
7	8	101	GLN
8	A	543	ASN
8	A	611	ASN
8	A	650	GLN
8	A	656	ASN
1	C	98	ASN
1	C	114	GLN
1	C	289	ASN

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Mol	Chain	Res	Type
1	C	315	GLN
1	C	352	HIS
8	D	543	ASN
8	D	611	ASN
8	D	650	GLN
8	D	656	ASN
2	H	60	ASN
3	L	79	GLN
4	M	23	ASN
4	M	60	ASN
7	Q	101	GLN
6	R	89	GLN
1	c	88	ASN
1	c	98	ASN
1	c	114	GLN
1	c	258	GLN
1	c	315	GLN
1	c	352	HIS
1	c	374	HIS
8	d	543	ASN
8	d	611	ASN
8	d	650	GLN
8	d	656	ASN
2	h	39	GLN
2	h	60	ASN
3	l	79	GLN
4	m	23	ASN
4	m	60	ASN
7	q	3	GLN
7	q	6	GLN
7	q	101	GLN
6	r	89	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

129 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
10	NAG	0	1	10,1	14,14,15	0.46	0	17,19,21	0.66	0
10	NAG	0	2	10	14,14,15	0.63	0	17,19,21	1.29	1 (5%)
10	BMA	0	3	10	11,11,12	0.64	0	15,15,17	1.34	2 (13%)
10	MAN	0	4	10	11,11,12	0.67	0	15,15,17	1.09	2 (13%)
13	NAG	1	1	13,1	14,14,15	0.50	0	17,19,21	0.53	0
13	NAG	1	2	13	14,14,15	0.24	0	17,19,21	1.21	2 (11%)
13	BMA	1	3	13	11,11,12	0.99	1 (9%)	15,15,17	1.30	2 (13%)
13	MAN	1	4	13	11,11,12	0.72	0	15,15,17	1.48	1 (6%)
13	MAN	1	5	13	11,11,12	1.09	1 (9%)	15,15,17	1.87	2 (13%)
13	MAN	1	6	13	11,11,12	0.93	0	15,15,17	0.93	2 (13%)
13	MAN	1	7	13	11,11,12	0.73	0	15,15,17	1.03	2 (13%)
13	MAN	1	8	13	11,11,12	0.68	0	15,15,17	1.18	2 (13%)
13	MAN	1	9	13	11,11,12	0.81	0	15,15,17	1.18	2 (13%)
9	NAG	9	1	9	14,14,15	0.75	1 (7%)	17,19,21	0.96	1 (5%)
9	NAG	9	2	9	14,14,15	0.23	0	17,19,21	0.65	1 (5%)
10	NAG	AA	1	10,1	14,14,15	0.24	0	17,19,21	0.72	0
10	NAG	AA	2	10	14,14,15	0.35	0	17,19,21	0.87	1 (5%)
10	BMA	AA	3	10	11,11,12	0.82	0	15,15,17	0.70	0
10	MAN	AA	4	10	11,11,12	0.95	1 (9%)	15,15,17	1.32	2 (13%)
9	NAG	B	1	9,1	14,14,15	0.51	0	17,19,21	0.96	1 (5%)
9	NAG	B	2	9	14,14,15	0.38	0	17,19,21	0.54	0
10	NAG	E	1	10,1	14,14,15	0.23	0	17,19,21	0.47	0
10	NAG	E	2	10	14,14,15	0.31	0	17,19,21	0.94	1 (5%)
10	BMA	E	3	10	11,11,12	0.70	0	15,15,17	1.08	1 (6%)
10	MAN	E	4	10	11,11,12	0.83	0	15,15,17	1.02	1 (6%)
9	NAG	F	1	9,1	14,14,15	0.35	0	17,19,21	0.48	0
9	NAG	F	2	9	14,14,15	0.26	0	17,19,21	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	NAG	G	1	9,1	14,14,15	0.52	0	17,19,21	1.17	2 (11%)
9	NAG	G	2	9	14,14,15	0.32	0	17,19,21	0.46	0
9	NAG	I	1	9,1	14,14,15	0.24	0	17,19,21	0.50	0
9	NAG	I	2	9	14,14,15	0.37	0	17,19,21	0.54	0
11	NAG	J	1	11,1	14,14,15	0.20	0	17,19,21	0.52	0
11	NAG	J	2	11	14,14,15	0.18	0	17,19,21	0.54	0
11	BMA	J	3	11	11,11,12	0.62	0	15,15,17	1.14	1 (6%)
11	MAN	J	4	11	11,11,12	0.73	0	15,15,17	0.97	2 (13%)
11	MAN	J	5	11	11,11,12	0.72	0	15,15,17	1.03	2 (13%)
9	NAG	K	1	9,1	14,14,15	0.35	0	17,19,21	0.58	0
9	NAG	K	2	9	14,14,15	0.36	0	17,19,21	0.53	0
9	NAG	O	1	9,1	14,14,15	0.23	0	17,19,21	0.69	1 (5%)
9	NAG	O	2	9	14,14,15	0.35	0	17,19,21	0.47	0
12	NAG	P	1	12,1	14,14,15	0.30	0	17,19,21	0.56	0
12	NAG	P	2	12	14,14,15	0.20	0	17,19,21	0.53	0
12	BMA	P	3	12	11,11,12	0.66	0	15,15,17	0.89	1 (6%)
10	NAG	S	1	10,1	14,14,15	0.38	0	17,19,21	0.65	0
10	NAG	S	2	10	14,14,15	0.63	0	17,19,21	1.30	1 (5%)
10	BMA	S	3	10	11,11,12	0.66	0	15,15,17	1.34	2 (13%)
10	MAN	S	4	10	11,11,12	0.67	0	15,15,17	1.08	2 (13%)
13	NAG	T	1	13,1	14,14,15	0.38	0	17,19,21	0.55	0
13	NAG	T	2	13	14,14,15	0.20	0	17,19,21	1.23	3 (17%)
13	BMA	T	3	13	11,11,12	1.05	1 (9%)	15,15,17	1.30	1 (6%)
13	MAN	T	4	13	11,11,12	0.63	0	15,15,17	1.37	1 (6%)
13	MAN	T	5	13	11,11,12	1.13	1 (9%)	15,15,17	1.82	2 (13%)
13	MAN	T	6	13	11,11,12	0.98	1 (9%)	15,15,17	0.95	2 (13%)
13	MAN	T	7	13	11,11,12	0.72	0	15,15,17	1.02	2 (13%)
13	MAN	T	8	13	11,11,12	0.72	0	15,15,17	1.17	2 (13%)
13	MAN	T	9	13	11,11,12	0.82	0	15,15,17	1.08	2 (13%)
9	NAG	U	1	9	14,14,15	0.74	1 (7%)	17,19,21	0.94	1 (5%)
9	NAG	U	2	9	14,14,15	0.25	0	17,19,21	0.65	1 (5%)
10	NAG	V	1	10,1	14,14,15	0.23	0	17,19,21	0.77	1 (5%)
10	NAG	V	2	10	14,14,15	0.38	0	17,19,21	0.91	1 (5%)
10	BMA	V	3	10	11,11,12	0.84	0	15,15,17	0.72	0
10	MAN	V	4	10	11,11,12	0.92	1 (9%)	15,15,17	1.28	2 (13%)
9	NAG	W	1	9,1	14,14,15	0.51	0	17,19,21	0.97	1 (5%)
9	NAG	W	2	9	14,14,15	0.39	0	17,19,21	0.55	0
10	NAG	X	1	10,1	14,14,15	0.26	0	17,19,21	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	NAG	X	2	10	14,14,15	0.26	0	17,19,21	0.93	1 (5%)
10	BMA	X	3	10	11,11,12	0.69	0	15,15,17	1.12	1 (6%)
10	MAN	X	4	10	11,11,12	0.85	0	15,15,17	1.00	1 (6%)
9	NAG	Y	1	9,1	14,14,15	0.31	0	17,19,21	0.50	0
9	NAG	Y	2	9	14,14,15	0.26	0	17,19,21	0.46	0
9	NAG	Z	1	9,1	14,14,15	0.54	0	17,19,21	1.12	2 (11%)
9	NAG	Z	2	9	14,14,15	0.24	0	17,19,21	0.46	0
9	NAG	a	1	9,1	14,14,15	0.23	0	17,19,21	0.52	0
9	NAG	a	2	9	14,14,15	0.36	0	17,19,21	0.52	0
11	NAG	b	1	11,1	14,14,15	0.21	0	17,19,21	0.47	0
11	NAG	b	2	11	14,14,15	0.18	0	17,19,21	0.55	0
11	BMA	b	3	11	11,11,12	0.69	0	15,15,17	1.10	1 (6%)
11	MAN	b	4	11	11,11,12	0.79	0	15,15,17	0.95	2 (13%)
11	MAN	b	5	11	11,11,12	0.74	0	15,15,17	1.01	2 (13%)
9	NAG	e	1	9,1	14,14,15	0.23	0	17,19,21	0.57	0
9	NAG	e	2	9	14,14,15	0.30	0	17,19,21	0.53	0
9	NAG	f	1	9,1	14,14,15	0.22	0	17,19,21	0.69	1 (5%)
9	NAG	f	2	9	14,14,15	0.38	0	17,19,21	0.47	0
12	NAG	g	1	12,1	14,14,15	0.41	0	17,19,21	0.64	1 (5%)
12	NAG	g	2	12	14,14,15	0.27	0	17,19,21	0.60	0
12	BMA	g	3	12	11,11,12	0.70	0	15,15,17	0.92	1 (6%)
10	NAG	i	1	10,1	14,14,15	0.32	0	17,19,21	0.71	1 (5%)
10	NAG	i	2	10	14,14,15	0.61	0	17,19,21	1.36	3 (17%)
10	BMA	i	3	10	11,11,12	0.69	0	15,15,17	1.42	2 (13%)
10	MAN	i	4	10	11,11,12	0.65	0	15,15,17	1.14	2 (13%)
13	NAG	j	1	13,1	14,14,15	0.41	0	17,19,21	0.59	0
13	NAG	j	2	13	14,14,15	0.20	0	17,19,21	1.21	3 (17%)
13	BMA	j	3	13	11,11,12	0.91	0	15,15,17	1.34	1 (6%)
13	MAN	j	4	13	11,11,12	0.64	0	15,15,17	1.33	1 (6%)
13	MAN	j	5	13	11,11,12	1.02	1 (9%)	15,15,17	1.77	2 (13%)
13	MAN	j	6	13	11,11,12	0.97	1 (9%)	15,15,17	0.98	2 (13%)
13	MAN	j	7	13	11,11,12	0.72	0	15,15,17	1.02	2 (13%)
13	MAN	j	8	13	11,11,12	0.73	0	15,15,17	1.18	2 (13%)
13	MAN	j	9	13	11,11,12	0.82	0	15,15,17	1.07	2 (13%)
9	NAG	k	1	9	14,14,15	0.79	1 (7%)	17,19,21	0.95	1 (5%)
9	NAG	k	2	9	14,14,15	0.27	0	17,19,21	0.66	1 (5%)
10	NAG	o	1	10,1	14,14,15	0.25	0	17,19,21	0.72	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	NAG	o	2	10	14,14,15	0.34	0	17,19,21	0.87	1 (5%)
10	BMA	o	3	10	11,11,12	0.80	0	15,15,17	0.70	0
10	MAN	o	4	10	11,11,12	0.92	1 (9%)	15,15,17	1.28	2 (13%)
9	NAG	p	1	9,1	14,14,15	0.44	0	17,19,21	0.97	1 (5%)
9	NAG	p	2	9	14,14,15	0.37	0	17,19,21	0.54	0
10	NAG	s	1	10,1	14,14,15	0.29	0	17,19,21	0.49	0
10	NAG	s	2	10	14,14,15	0.30	0	17,19,21	0.91	1 (5%)
10	BMA	s	3	10	11,11,12	0.72	0	15,15,17	1.02	1 (6%)
10	MAN	s	4	10	11,11,12	0.83	0	15,15,17	1.05	1 (6%)
9	NAG	t	1	9,1	14,14,15	0.41	0	17,19,21	0.49	0
9	NAG	t	2	9	14,14,15	0.25	0	17,19,21	0.52	0
9	NAG	u	1	9,1	14,14,15	0.46	0	17,19,21	1.16	2 (11%)
9	NAG	u	2	9	14,14,15	0.29	0	17,19,21	0.47	0
9	NAG	v	1	9,1	14,14,15	0.28	0	17,19,21	0.58	0
9	NAG	v	2	9	14,14,15	0.30	0	17,19,21	0.53	0
11	NAG	w	1	11,1	14,14,15	0.19	0	17,19,21	0.55	0
11	NAG	w	2	11	14,14,15	0.18	0	17,19,21	0.49	0
11	BMA	w	3	11	11,11,12	0.58	0	15,15,17	1.10	1 (6%)
11	MAN	w	4	11	11,11,12	0.76	0	15,15,17	0.96	2 (13%)
11	MAN	w	5	11	11,11,12	0.74	0	15,15,17	1.03	2 (13%)
9	NAG	x	1	9,1	14,14,15	0.37	0	17,19,21	0.61	0
9	NAG	x	2	9	14,14,15	0.41	0	17,19,21	0.50	0
9	NAG	y	1	9,1	14,14,15	0.22	0	17,19,21	0.57	0
9	NAG	y	2	9	14,14,15	0.30	0	17,19,21	0.48	0
12	NAG	z	1	12,1	14,14,15	0.35	0	17,19,21	0.48	0
12	NAG	z	2	12	14,14,15	0.20	0	17,19,21	0.54	0
12	BMA	z	3	12	11,11,12	0.67	0	15,15,17	0.93	2 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	NAG	0	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	0	2	10	-	2/6/23/26	0/1/1/1
10	BMA	0	3	10	-	2/2/19/22	0/1/1/1
10	MAN	0	4	10	-	0/2/19/22	0/1/1/1
13	NAG	1	1	13,1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
13	NAG	1	2	13	-	4/6/23/26	0/1/1/1
13	BMA	1	3	13	-	0/2/19/22	0/1/1/1
13	MAN	1	4	13	-	0/2/19/22	0/1/1/1
13	MAN	1	5	13	-	2/2/19/22	0/1/1/1
13	MAN	1	6	13	-	0/2/19/22	0/1/1/1
13	MAN	1	7	13	-	2/2/19/22	0/1/1/1
13	MAN	1	8	13	-	2/2/19/22	0/1/1/1
13	MAN	1	9	13	-	2/2/19/22	0/1/1/1
9	NAG	9	1	9	-	2/6/23/26	0/1/1/1
9	NAG	9	2	9	-	2/6/23/26	0/1/1/1
10	NAG	AA	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	AA	2	10	-	0/6/23/26	0/1/1/1
10	BMA	AA	3	10	-	0/2/19/22	0/1/1/1
10	MAN	AA	4	10	-	2/2/19/22	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
10	NAG	E	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	E	2	10	-	2/6/23/26	0/1/1/1
10	BMA	E	3	10	-	2/2/19/22	0/1/1/1
10	MAN	E	4	10	-	0/2/19/22	0/1/1/1
9	NAG	F	1	9,1	-	1/6/23/26	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
9	NAG	G	2	9	-	0/6/23/26	0/1/1/1
9	NAG	I	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	I	2	9	-	0/6/23/26	0/1/1/1
11	NAG	J	1	11,1	-	2/6/23/26	0/1/1/1
11	NAG	J	2	11	-	0/6/23/26	0/1/1/1
11	BMA	J	3	11	-	1/2/19/22	0/1/1/1
11	MAN	J	4	11	-	1/2/19/22	0/1/1/1
11	MAN	J	5	11	-	0/2/19/22	0/1/1/1
9	NAG	K	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	K	2	9	-	0/6/23/26	0/1/1/1
9	NAG	O	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	O	2	9	-	2/6/23/26	0/1/1/1
12	NAG	P	1	12,1	-	2/6/23/26	0/1/1/1
12	NAG	P	2	12	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	BMA	P	3	12	-	2/2/19/22	0/1/1/1
10	NAG	S	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	S	2	10	-	2/6/23/26	0/1/1/1
10	BMA	S	3	10	-	2/2/19/22	0/1/1/1
10	MAN	S	4	10	-	0/2/19/22	0/1/1/1
13	NAG	T	1	13,1	-	2/6/23/26	0/1/1/1
13	NAG	T	2	13	-	4/6/23/26	0/1/1/1
13	BMA	T	3	13	-	0/2/19/22	0/1/1/1
13	MAN	T	4	13	-	0/2/19/22	0/1/1/1
13	MAN	T	5	13	-	2/2/19/22	0/1/1/1
13	MAN	T	6	13	-	0/2/19/22	0/1/1/1
13	MAN	T	7	13	-	2/2/19/22	0/1/1/1
13	MAN	T	8	13	-	2/2/19/22	0/1/1/1
13	MAN	T	9	13	-	2/2/19/22	0/1/1/1
9	NAG	U	1	9	-	3/6/23/26	0/1/1/1
9	NAG	U	2	9	-	2/6/23/26	0/1/1/1
10	NAG	V	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	V	2	10	-	0/6/23/26	0/1/1/1
10	BMA	V	3	10	-	0/2/19/22	0/1/1/1
10	MAN	V	4	10	-	2/2/19/22	0/1/1/1
9	NAG	W	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	W	2	9	-	2/6/23/26	0/1/1/1
10	NAG	X	1	10,1	-	2/6/23/26	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
10	MAN	X	4	10	-	0/2/19/22	0/1/1/1
9	NAG	Y	1	9,1	-	2/6/23/26	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
9	NAG	Z	2	9	-	0/6/23/26	0/1/1/1
9	NAG	a	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	a	2	9	-	0/6/23/26	0/1/1/1
11	NAG	b	1	11,1	-	2/6/23/26	0/1/1/1
11	NAG	b	2	11	-	0/6/23/26	0/1/1/1
11	BMA	b	3	11	-	1/2/19/22	0/1/1/1
11	MAN	b	4	11	-	1/2/19/22	0/1/1/1
11	MAN	b	5	11	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	e	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	e	2	9	-	0/6/23/26	0/1/1/1
9	NAG	f	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	f	2	9	-	2/6/23/26	0/1/1/1
12	NAG	g	1	12,1	-	2/6/23/26	0/1/1/1
12	NAG	g	2	12	-	0/6/23/26	0/1/1/1
12	BMA	g	3	12	-	2/2/19/22	0/1/1/1
10	NAG	i	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	i	2	10	-	2/6/23/26	0/1/1/1
10	BMA	i	3	10	-	2/2/19/22	0/1/1/1
10	MAN	i	4	10	-	0/2/19/22	0/1/1/1
13	NAG	j	1	13,1	-	2/6/23/26	0/1/1/1
13	NAG	j	2	13	-	4/6/23/26	0/1/1/1
13	BMA	j	3	13	-	0/2/19/22	0/1/1/1
13	MAN	j	4	13	-	2/2/19/22	0/1/1/1
13	MAN	j	5	13	-	2/2/19/22	0/1/1/1
13	MAN	j	6	13	-	0/2/19/22	0/1/1/1
13	MAN	j	7	13	-	2/2/19/22	0/1/1/1
13	MAN	j	8	13	-	2/2/19/22	0/1/1/1
13	MAN	j	9	13	-	2/2/19/22	0/1/1/1
9	NAG	k	1	9	-	2/6/23/26	0/1/1/1
9	NAG	k	2	9	-	2/6/23/26	0/1/1/1
10	NAG	o	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	o	2	10	-	0/6/23/26	0/1/1/1
10	BMA	o	3	10	-	0/2/19/22	0/1/1/1
10	MAN	o	4	10	-	2/2/19/22	0/1/1/1
9	NAG	p	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	p	2	9	-	2/6/23/26	0/1/1/1
10	NAG	s	1	10,1	-	2/6/23/26	0/1/1/1
10	NAG	s	2	10	-	2/6/23/26	0/1/1/1
10	BMA	s	3	10	-	2/2/19/22	0/1/1/1
10	MAN	s	4	10	-	0/2/19/22	0/1/1/1
9	NAG	t	1	9,1	-	1/6/23/26	0/1/1/1
9	NAG	t	2	9	-	0/6/23/26	0/1/1/1
9	NAG	u	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	u	2	9	-	0/6/23/26	0/1/1/1
9	NAG	v	1	9,1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	NAG	v	2	9	-	0/6/23/26	0/1/1/1
11	NAG	w	1	11,1	-	1/6/23/26	0/1/1/1
11	NAG	w	2	11	-	0/6/23/26	0/1/1/1
11	BMA	w	3	11	-	1/2/19/22	0/1/1/1
11	MAN	w	4	11	-	1/2/19/22	0/1/1/1
11	MAN	w	5	11	-	0/2/19/22	0/1/1/1
9	NAG	x	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	x	2	9	-	0/6/23/26	0/1/1/1
9	NAG	y	1	9,1	-	2/6/23/26	0/1/1/1
9	NAG	y	2	9	-	2/6/23/26	0/1/1/1
12	NAG	z	1	12,1	-	2/6/23/26	0/1/1/1
12	NAG	z	2	12	-	0/6/23/26	0/1/1/1
12	BMA	z	3	12	-	2/2/19/22	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	AA	4	MAN	C1-C2	2.74	1.58	1.52
10	V	4	MAN	C1-C2	2.62	1.58	1.52
10	o	4	MAN	C1-C2	2.60	1.58	1.52
9	k	1	NAG	C1-C2	2.52	1.55	1.52
9	U	1	NAG	C1-C2	2.42	1.55	1.52
9	9	1	NAG	C1-C2	2.29	1.55	1.52
13	1	5	MAN	O5-C5	2.28	1.47	1.43
13	T	3	BMA	O5-C1	-2.22	1.40	1.43
13	T	5	MAN	O5-C5	2.22	1.47	1.43
13	1	3	BMA	O5-C1	-2.20	1.40	1.43
13	j	5	MAN	O5-C5	2.20	1.47	1.43
13	T	6	MAN	O5-C1	-2.11	1.40	1.43
13	j	6	MAN	O5-C1	-2.03	1.40	1.43

All (119) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	1	5	MAN	C1-O5-C5	5.57	119.65	112.19
13	T	5	MAN	C1-O5-C5	5.22	119.18	112.19
13	1	4	MAN	C1-O5-C5	5.19	119.14	112.19
13	j	5	MAN	C1-O5-C5	5.16	119.10	112.19
13	T	4	MAN	C1-O5-C5	4.66	118.44	112.19
10	i	3	BMA	C1-O5-C5	4.44	118.14	112.19
13	j	4	MAN	C1-O5-C5	4.42	118.11	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	0	3	BMA	C1-O5-C5	4.25	117.88	112.19
10	S	3	BMA	C1-O5-C5	4.17	117.77	112.19
10	i	2	NAG	C1-O5-C5	4.12	117.71	112.19
13	T	5	MAN	O2-C2-C3	-3.97	101.94	110.15
10	S	2	NAG	C1-O5-C5	3.95	117.48	112.19
13	1	5	MAN	O2-C2-C3	-3.95	101.97	110.15
10	0	2	NAG	C1-O5-C5	3.91	117.43	112.19
13	j	3	BMA	C1-O5-C5	3.80	117.28	112.19
13	j	5	MAN	O2-C2-C3	-3.78	102.31	110.15
10	AA	4	MAN	C1-O5-C5	3.67	117.10	112.19
13	T	3	BMA	C1-O5-C5	3.55	116.94	112.19
13	1	3	BMA	C1-O5-C5	3.55	116.94	112.19
10	X	3	BMA	C1-O5-C5	3.53	116.92	112.19
10	o	4	MAN	C1-O5-C5	3.52	116.90	112.19
10	V	4	MAN	C1-O5-C5	3.50	116.88	112.19
13	1	9	MAN	C1-O5-C5	3.45	116.81	112.19
13	j	8	MAN	C1-O5-C5	3.34	116.67	112.19
10	E	3	BMA	C1-O5-C5	3.33	116.65	112.19
13	1	8	MAN	C1-O5-C5	3.33	116.64	112.19
13	T	8	MAN	C1-O5-C5	3.31	116.62	112.19
10	i	4	MAN	C1-O5-C5	3.28	116.58	112.19
13	T	2	NAG	C2-N2-C7	3.26	127.27	122.90
13	1	2	NAG	C2-N2-C7	3.26	127.27	122.90
9	W	1	NAG	C2-N2-C7	3.25	127.25	122.90
13	j	2	NAG	C2-N2-C7	3.24	127.25	122.90
9	B	1	NAG	C2-N2-C7	3.23	127.23	122.90
9	p	1	NAG	C2-N2-C7	3.20	127.19	122.90
10	s	3	BMA	C1-O5-C5	3.17	116.43	112.19
9	9	1	NAG	C2-N2-C7	3.14	127.10	122.90
10	s	4	MAN	C1-O5-C5	3.07	116.30	112.19
9	k	1	NAG	C2-N2-C7	3.06	127.01	122.90
10	0	4	MAN	C1-O5-C5	3.06	116.29	112.19
9	U	1	NAG	C2-N2-C7	3.05	126.98	122.90
10	S	4	MAN	C1-O5-C5	3.04	116.25	112.19
13	T	9	MAN	C1-O5-C5	3.02	116.24	112.19
11	J	3	BMA	C1-O5-C5	3.01	116.23	112.19
9	Z	1	NAG	C2-N2-C7	3.00	126.92	122.90
9	G	1	NAG	C2-N2-C7	3.00	126.92	122.90
9	u	1	NAG	C2-N2-C7	2.99	126.90	122.90
10	E	4	MAN	C1-O5-C5	2.94	116.12	112.19
10	E	2	NAG	C1-O5-C5	2.92	116.10	112.19
13	j	9	MAN	C1-O5-C5	2.90	116.08	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	w	3	BMA	C1-O5-C5	2.90	116.08	112.19
10	X	2	NAG	C1-O5-C5	2.88	116.04	112.19
10	s	2	NAG	C1-O5-C5	2.86	116.02	112.19
10	X	4	MAN	C1-O5-C5	2.77	115.89	112.19
11	w	5	MAN	C1-O5-C5	2.67	115.76	112.19
11	J	5	MAN	C1-O5-C5	2.65	115.74	112.19
10	V	2	NAG	C1-O5-C5	2.63	115.71	112.19
9	G	1	NAG	C1-O5-C5	2.59	115.66	112.19
10	o	2	NAG	C1-O5-C5	2.59	115.66	112.19
13	T	7	MAN	O2-C2-C3	-2.57	104.83	110.15
12	z	3	BMA	C1-O5-C5	2.56	115.62	112.19
13	1	7	MAN	O2-C2-C3	-2.55	104.86	110.15
13	j	7	MAN	O2-C2-C3	-2.55	104.88	110.15
9	u	1	NAG	C1-O5-C5	2.53	115.58	112.19
12	g	3	BMA	C1-O5-C5	2.51	115.55	112.19
11	b	3	BMA	C1-O5-C5	2.48	115.50	112.19
11	b	5	MAN	C1-O5-C5	2.47	115.49	112.19
10	AA	2	NAG	C1-O5-C5	2.45	115.47	112.19
13	j	7	MAN	C1-O5-C5	2.45	115.47	112.19
13	T	7	MAN	C1-O5-C5	2.44	115.45	112.19
13	1	7	MAN	C1-O5-C5	2.42	115.43	112.19
11	b	5	MAN	O2-C2-C3	-2.39	105.20	110.15
13	j	8	MAN	O2-C2-C3	-2.37	105.24	110.15
13	T	8	MAN	O2-C2-C3	-2.37	105.25	110.15
13	j	6	MAN	C1-O5-C5	2.35	115.33	112.19
13	T	6	MAN	O2-C2-C3	-2.35	105.29	110.15
13	j	6	MAN	O2-C2-C3	-2.34	105.31	110.15
10	i	4	MAN	O2-C2-C3	-2.33	105.32	110.15
10	0	4	MAN	O2-C2-C3	-2.33	105.32	110.15
11	J	5	MAN	O2-C2-C3	-2.33	105.33	110.15
13	1	8	MAN	O2-C2-C3	-2.32	105.34	110.15
12	P	3	BMA	C1-O5-C5	2.32	115.30	112.19
10	i	1	NAG	C1-O5-C5	2.30	115.27	112.19
9	O	1	NAG	C1-O5-C5	2.30	115.27	112.19
9	f	1	NAG	C1-O5-C5	2.30	115.27	112.19
11	w	5	MAN	O2-C2-C3	-2.30	105.39	110.15
11	J	4	MAN	O2-C2-C3	-2.29	105.41	110.15
11	b	4	MAN	O2-C2-C3	-2.29	105.41	110.15
10	AA	4	MAN	O2-C2-C3	-2.27	105.44	110.15
9	Z	1	NAG	C1-O5-C5	2.27	115.23	112.19
10	o	4	MAN	O2-C2-C3	-2.26	105.47	110.15
10	S	4	MAN	O2-C2-C3	-2.26	105.47	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	V	4	MAN	O2-C2-C3	-2.25	105.48	110.15
11	w	4	MAN	O2-C2-C3	-2.25	105.49	110.15
13	1	6	MAN	O2-C2-C3	-2.22	105.55	110.15
9	9	2	NAG	C1-O5-C5	2.21	115.15	112.19
9	k	2	NAG	C1-O5-C5	2.20	115.14	112.19
13	T	2	NAG	C1-C2-N2	2.20	113.90	110.43
11	w	4	MAN	C1-O5-C5	2.20	115.13	112.19
11	J	4	MAN	C1-O5-C5	2.19	115.12	112.19
9	U	2	NAG	C1-O5-C5	2.17	115.09	112.19
13	T	6	MAN	C1-O5-C5	2.16	115.08	112.19
13	j	2	NAG	C1-C2-N2	2.16	113.83	110.43
10	V	1	NAG	C1-O5-C5	2.14	115.05	112.19
13	1	6	MAN	C1-O5-C5	2.11	115.02	112.19
11	b	4	MAN	C1-O5-C5	2.10	115.00	112.19
13	1	2	NAG	C1-C2-N2	2.10	113.74	110.43
10	i	3	BMA	O2-C2-C3	-2.10	105.80	110.15
12	z	3	BMA	O2-C2-C3	-2.08	105.85	110.15
10	i	2	NAG	O4-C4-C3	2.06	115.23	110.38
13	j	2	NAG	C1-O5-C5	2.05	114.94	112.19
13	1	9	MAN	O2-C2-C3	-2.05	105.90	110.15
13	j	9	MAN	O2-C2-C3	-2.05	105.91	110.15
13	T	9	MAN	O2-C2-C3	-2.04	105.92	110.15
13	T	2	NAG	C1-O5-C5	2.04	114.92	112.19
13	1	3	BMA	O2-C2-C3	-2.03	105.94	110.15
10	S	3	BMA	O2-C2-C3	-2.01	105.98	110.15
10	0	3	BMA	O2-C2-C3	-2.01	105.98	110.15
10	i	2	NAG	O4-C4-C5	2.01	114.28	109.32
12	g	1	NAG	C1-O5-C5	2.01	114.87	112.19

There are no chirality outliers.

All (166) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	i	2	NAG	O5-C5-C6-O6
13	T	1	NAG	O5-C5-C6-O6
9	p	2	NAG	O5-C5-C6-O6
10	S	2	NAG	O5-C5-C6-O6
9	W	2	NAG	O5-C5-C6-O6
9	y	2	NAG	O5-C5-C6-O6
10	s	2	NAG	O5-C5-C6-O6
10	0	2	NAG	O5-C5-C6-O6
13	1	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
10	E	2	NAG	O5-C5-C6-O6
10	o	1	NAG	C4-C5-C6-O6
12	g	3	BMA	C4-C5-C6-O6
12	z	1	NAG	C4-C5-C6-O6
9	O	2	NAG	O5-C5-C6-O6
9	f	2	NAG	O5-C5-C6-O6
10	X	2	NAG	O5-C5-C6-O6
10	AA	1	NAG	C4-C5-C6-O6
12	P	1	NAG	C4-C5-C6-O6
12	z	3	BMA	C4-C5-C6-O6
9	B	2	NAG	O5-C5-C6-O6
13	T	5	MAN	O5-C5-C6-O6
13	j	1	NAG	O5-C5-C6-O6
10	V	1	NAG	C4-C5-C6-O6
12	P	3	BMA	C4-C5-C6-O6
13	1	5	MAN	C4-C5-C6-O6
10	i	1	NAG	O5-C5-C6-O6
10	0	3	BMA	O5-C5-C6-O6
13	j	5	MAN	O5-C5-C6-O6
9	y	2	NAG	C4-C5-C6-O6
10	s	2	NAG	C4-C5-C6-O6
13	1	1	NAG	C4-C5-C6-O6
10	i	3	BMA	O5-C5-C6-O6
12	z	1	NAG	O5-C5-C6-O6
13	T	8	MAN	O5-C5-C6-O6
13	j	8	MAN	O5-C5-C6-O6
10	E	2	NAG	C4-C5-C6-O6
10	X	2	NAG	C4-C5-C6-O6
10	0	2	NAG	C4-C5-C6-O6
12	g	1	NAG	C4-C5-C6-O6
13	T	1	NAG	C4-C5-C6-O6
10	S	3	BMA	O5-C5-C6-O6
13	j	2	NAG	O5-C5-C6-O6
13	1	8	MAN	O5-C5-C6-O6
9	U	2	NAG	O5-C5-C6-O6
9	k	2	NAG	O5-C5-C6-O6
9	W	2	NAG	C4-C5-C6-O6
10	i	2	NAG	C4-C5-C6-O6
10	o	1	NAG	O5-C5-C6-O6
9	p	2	NAG	C4-C5-C6-O6
10	S	2	NAG	C4-C5-C6-O6
10	i	1	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
10	AA	1	NAG	O5-C5-C6-O6
10	o	4	MAN	O5-C5-C6-O6
12	P	1	NAG	O5-C5-C6-O6
12	P	3	BMA	O5-C5-C6-O6
12	g	3	BMA	O5-C5-C6-O6
12	z	3	BMA	O5-C5-C6-O6
13	1	5	MAN	O5-C5-C6-O6
10	X	1	NAG	C4-C5-C6-O6
13	j	2	NAG	C4-C5-C6-O6
9	9	2	NAG	O5-C5-C6-O6
9	O	2	NAG	C4-C5-C6-O6
9	f	2	NAG	C4-C5-C6-O6
10	S	1	NAG	O5-C5-C6-O6
10	AA	4	MAN	O5-C5-C6-O6
9	B	2	NAG	C4-C5-C6-O6
13	j	1	NAG	C4-C5-C6-O6
13	T	2	NAG	O5-C5-C6-O6
13	1	2	NAG	O5-C5-C6-O6
10	o	4	MAN	C4-C5-C6-O6
10	V	1	NAG	O5-C5-C6-O6
10	X	1	NAG	O5-C5-C6-O6
10	E	1	NAG	C4-C5-C6-O6
10	V	4	MAN	O5-C5-C6-O6
10	X	3	BMA	C4-C5-C6-O6
13	T	2	NAG	C4-C5-C6-O6
13	T	5	MAN	C4-C5-C6-O6
13	1	2	NAG	C4-C5-C6-O6
12	g	1	NAG	O5-C5-C6-O6
9	k	2	NAG	C4-C5-C6-O6
9	U	2	NAG	C4-C5-C6-O6
10	s	1	NAG	C4-C5-C6-O6
10	V	4	MAN	C4-C5-C6-O6
10	AA	4	MAN	C4-C5-C6-O6
11	b	4	MAN	O5-C5-C6-O6
9	y	1	NAG	C4-C5-C6-O6
10	s	3	BMA	C4-C5-C6-O6
13	j	5	MAN	C4-C5-C6-O6
10	E	1	NAG	O5-C5-C6-O6
13	1	8	MAN	C4-C5-C6-O6
9	9	2	NAG	C4-C5-C6-O6
10	S	1	NAG	C4-C5-C6-O6
11	w	4	MAN	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
10	0	3	BMA	C4-C5-C6-O6
10	X	3	BMA	O5-C5-C6-O6
10	0	1	NAG	O5-C5-C6-O6
13	1	7	MAN	O5-C5-C6-O6
13	j	8	MAN	C4-C5-C6-O6
13	1	7	MAN	C4-C5-C6-O6
10	i	3	BMA	C4-C5-C6-O6
10	s	1	NAG	O5-C5-C6-O6
13	T	7	MAN	C4-C5-C6-O6
10	E	3	BMA	C4-C5-C6-O6
10	S	3	BMA	C4-C5-C6-O6
11	J	4	MAN	O5-C5-C6-O6
9	y	1	NAG	O5-C5-C6-O6
13	T	7	MAN	O5-C5-C6-O6
13	T	8	MAN	C4-C5-C6-O6
10	s	3	BMA	O5-C5-C6-O6
13	j	7	MAN	C4-C5-C6-O6
9	O	1	NAG	C4-C5-C6-O6
9	F	2	NAG	C4-C5-C6-O6
13	j	7	MAN	O5-C5-C6-O6
9	f	1	NAG	C4-C5-C6-O6
13	1	9	MAN	C4-C5-C6-O6
13	j	9	MAN	C4-C5-C6-O6
13	j	4	MAN	C4-C5-C6-O6
13	j	4	MAN	O5-C5-C6-O6
11	b	1	NAG	C4-C5-C6-O6
11	J	1	NAG	C4-C5-C6-O6
10	E	3	BMA	O5-C5-C6-O6
9	Y	2	NAG	C4-C5-C6-O6
9	F	2	NAG	O5-C5-C6-O6
13	T	9	MAN	C4-C5-C6-O6
9	Y	1	NAG	C4-C5-C6-O6
9	O	1	NAG	O5-C5-C6-O6
11	w	3	BMA	C4-C5-C6-O6
10	0	1	NAG	C4-C5-C6-O6
9	F	1	NAG	C4-C5-C6-O6
9	G	1	NAG	C3-C2-N2-C7
9	U	1	NAG	C3-C2-N2-C7
9	Z	1	NAG	C3-C2-N2-C7
9	k	1	NAG	C3-C2-N2-C7
9	u	1	NAG	C3-C2-N2-C7
9	9	1	NAG	C3-C2-N2-C7

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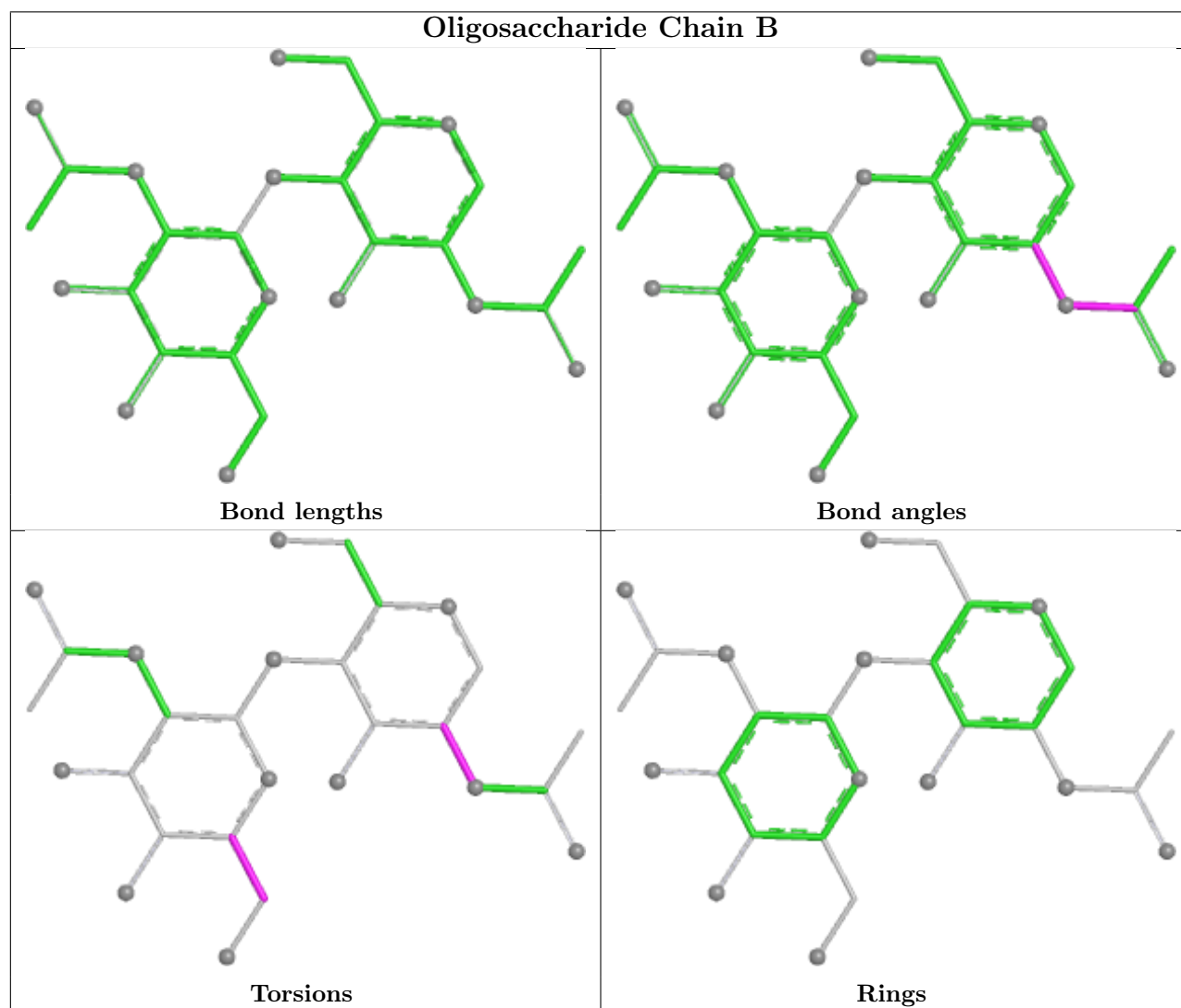
Mol	Chain	Res	Type	Atoms
13	j	2	NAG	C3-C2-N2-C7
13	l	2	NAG	C3-C2-N2-C7
13	j	9	MAN	O5-C5-C6-O6
13	l	9	MAN	O5-C5-C6-O6
9	f	1	NAG	O5-C5-C6-O6
9	U	1	NAG	O5-C5-C6-O6
11	J	3	BMA	C4-C5-C6-O6
11	b	1	NAG	O5-C5-C6-O6
11	w	1	NAG	C4-C5-C6-O6
9	Y	2	NAG	O5-C5-C6-O6
11	J	1	NAG	O5-C5-C6-O6
9	B	1	NAG	C1-C2-N2-C7
9	G	1	NAG	C1-C2-N2-C7
9	U	1	NAG	C1-C2-N2-C7
9	W	1	NAG	C1-C2-N2-C7
9	Z	1	NAG	C1-C2-N2-C7
9	k	1	NAG	C1-C2-N2-C7
9	p	1	NAG	C1-C2-N2-C7
9	u	1	NAG	C1-C2-N2-C7
9	9	1	NAG	C1-C2-N2-C7
13	T	2	NAG	C1-C2-N2-C7
13	j	2	NAG	C1-C2-N2-C7
13	l	2	NAG	C1-C2-N2-C7
13	T	9	MAN	O5-C5-C6-O6
9	B	1	NAG	C3-C2-N2-C7
9	W	1	NAG	C3-C2-N2-C7
9	p	1	NAG	C3-C2-N2-C7
13	T	2	NAG	C3-C2-N2-C7
11	b	3	BMA	C4-C5-C6-O6
9	t	1	NAG	C4-C5-C6-O6
9	Y	1	NAG	O5-C5-C6-O6

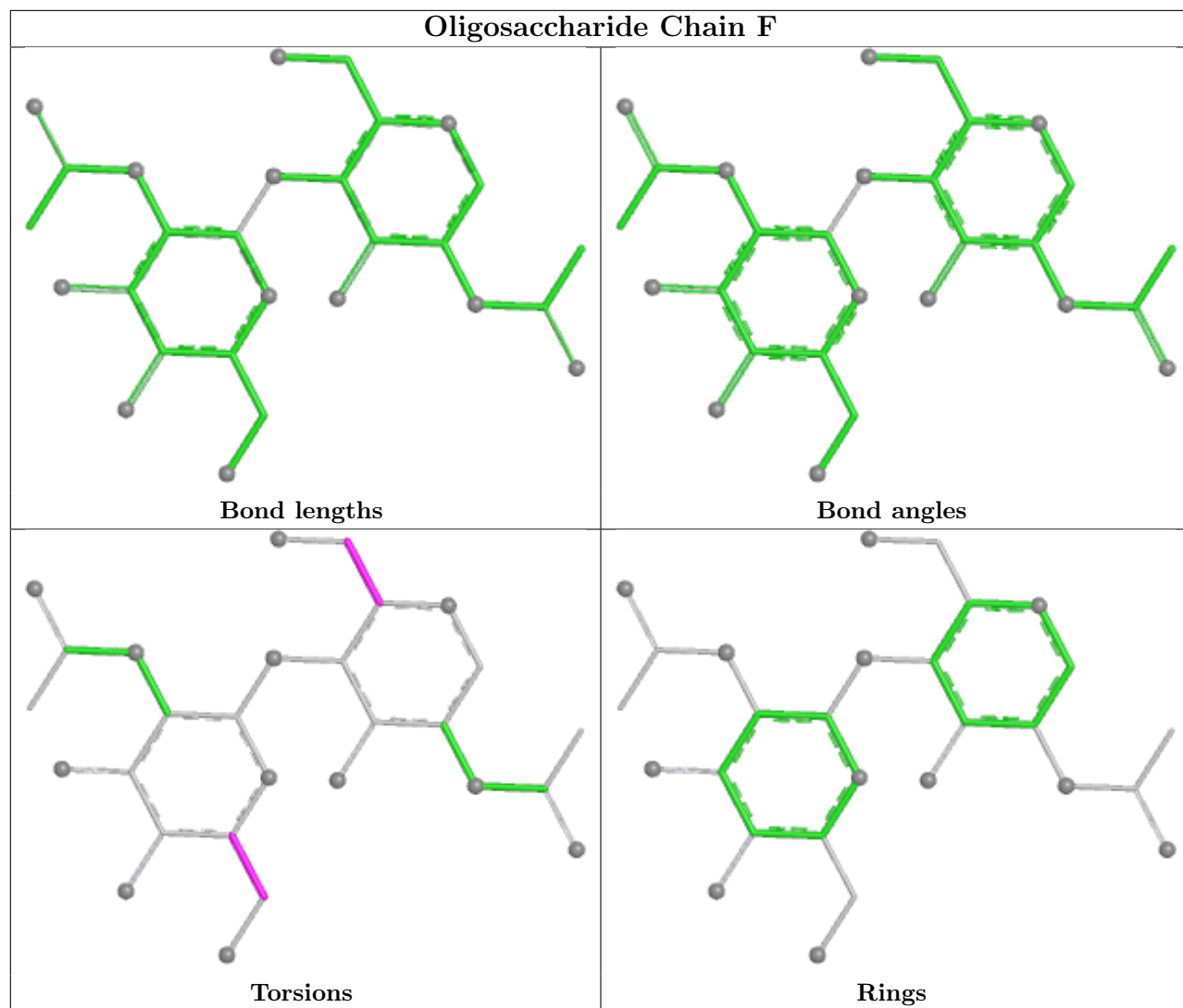
There are no ring outliers.

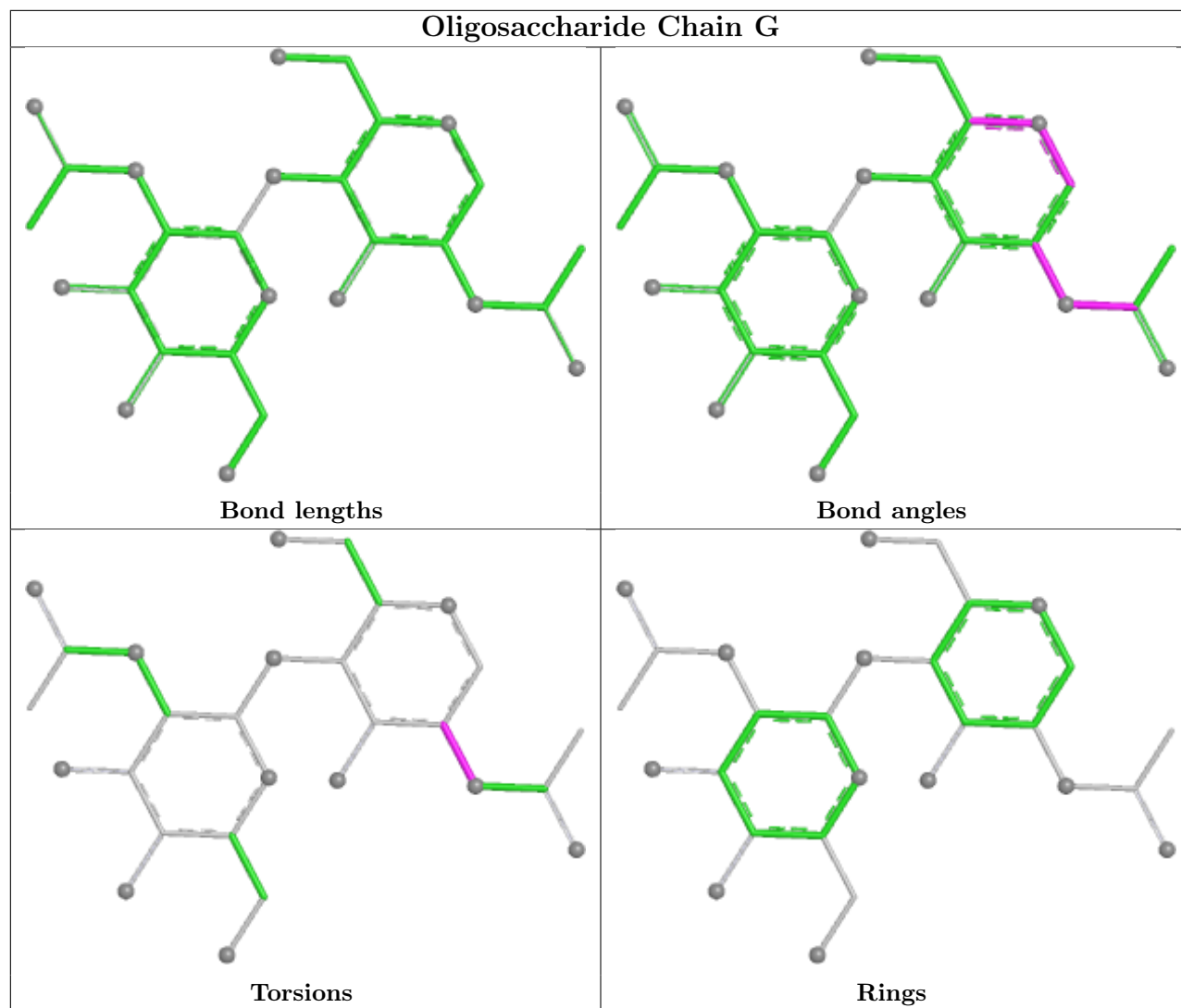
5 monomers are involved in 6 short contacts:

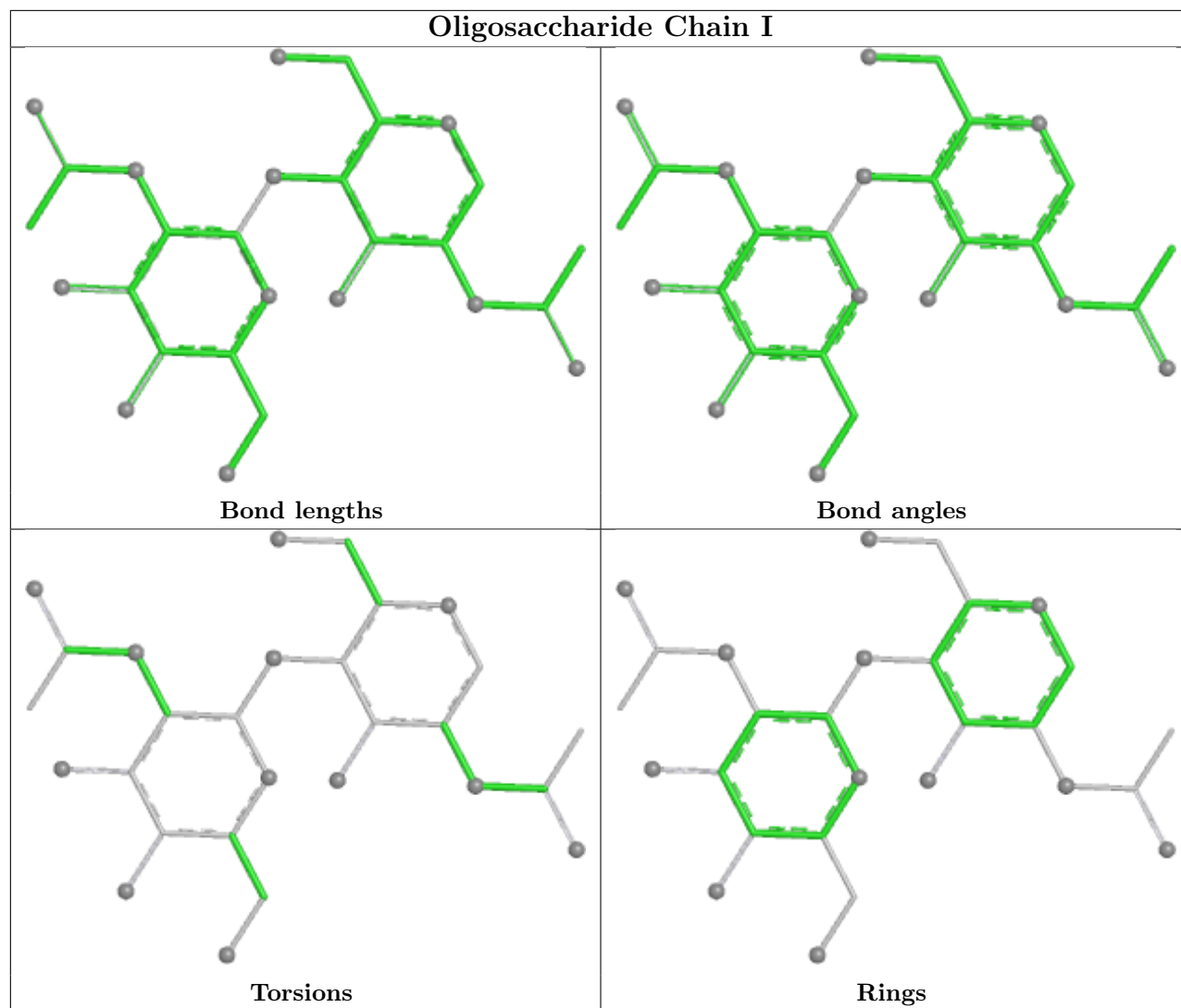
Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	T	5	MAN	1	0
10	V	2	NAG	1	0
10	AA	2	NAG	1	0
10	o	2	NAG	1	0
13	j	5	MAN	2	0

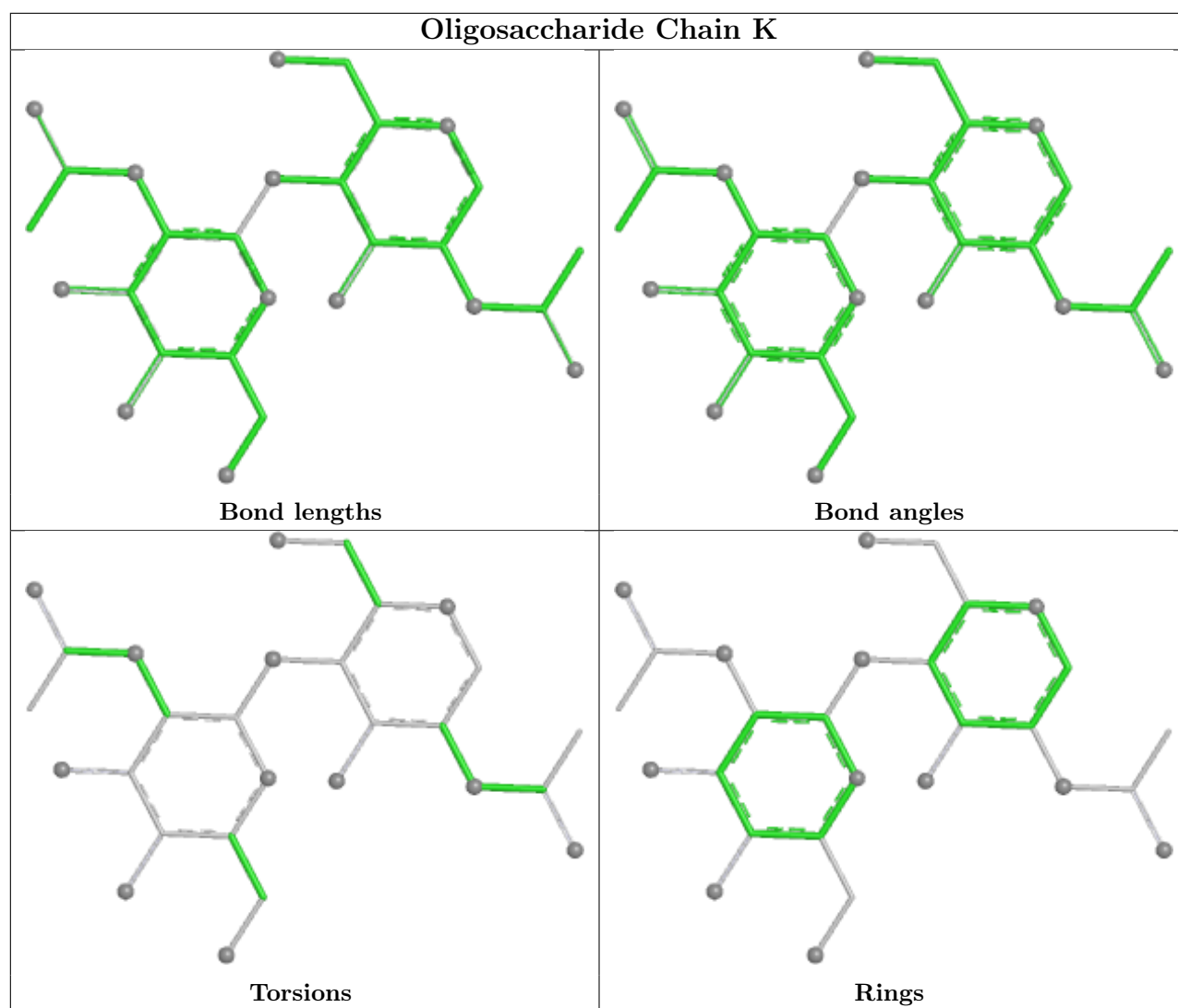
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

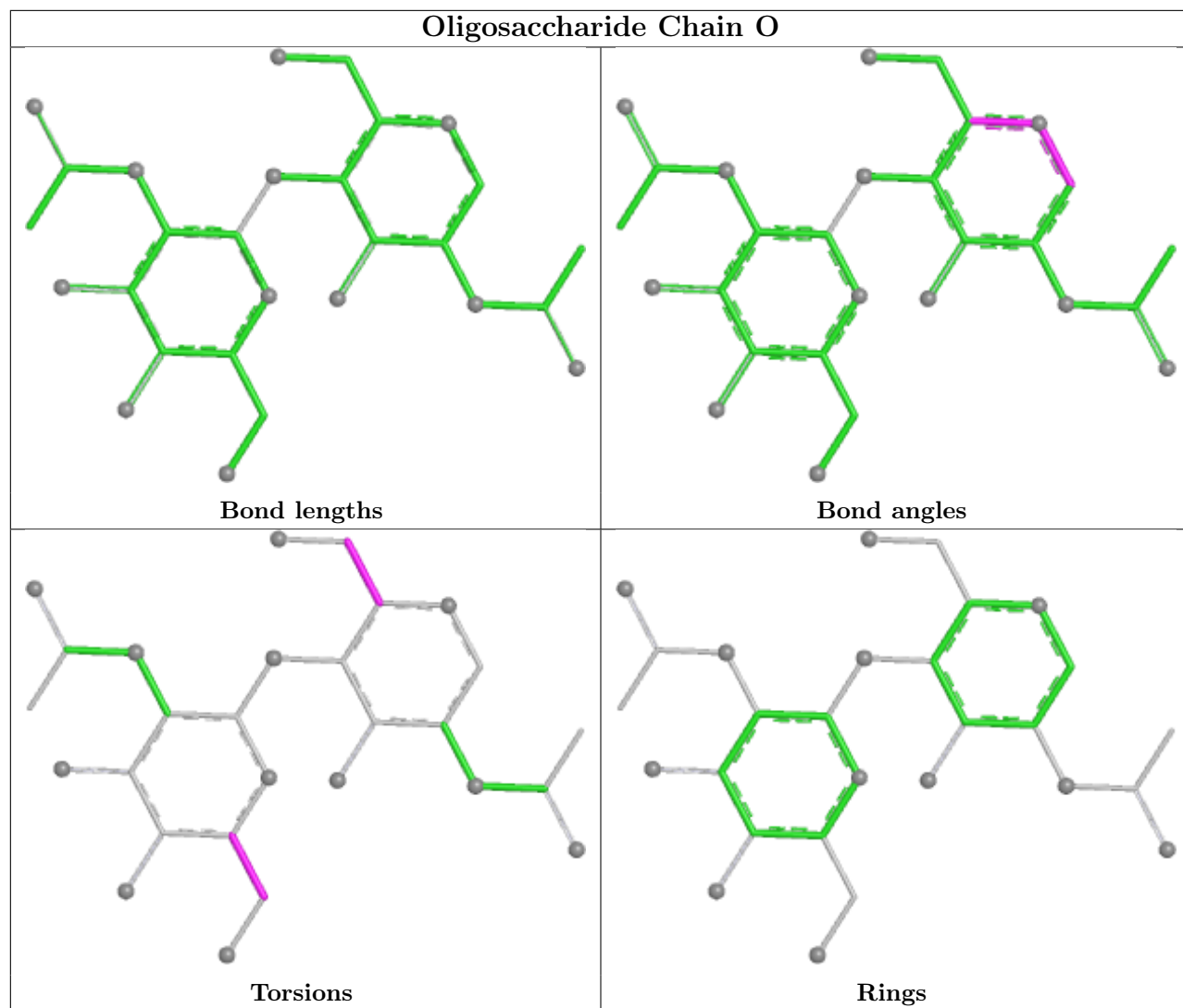


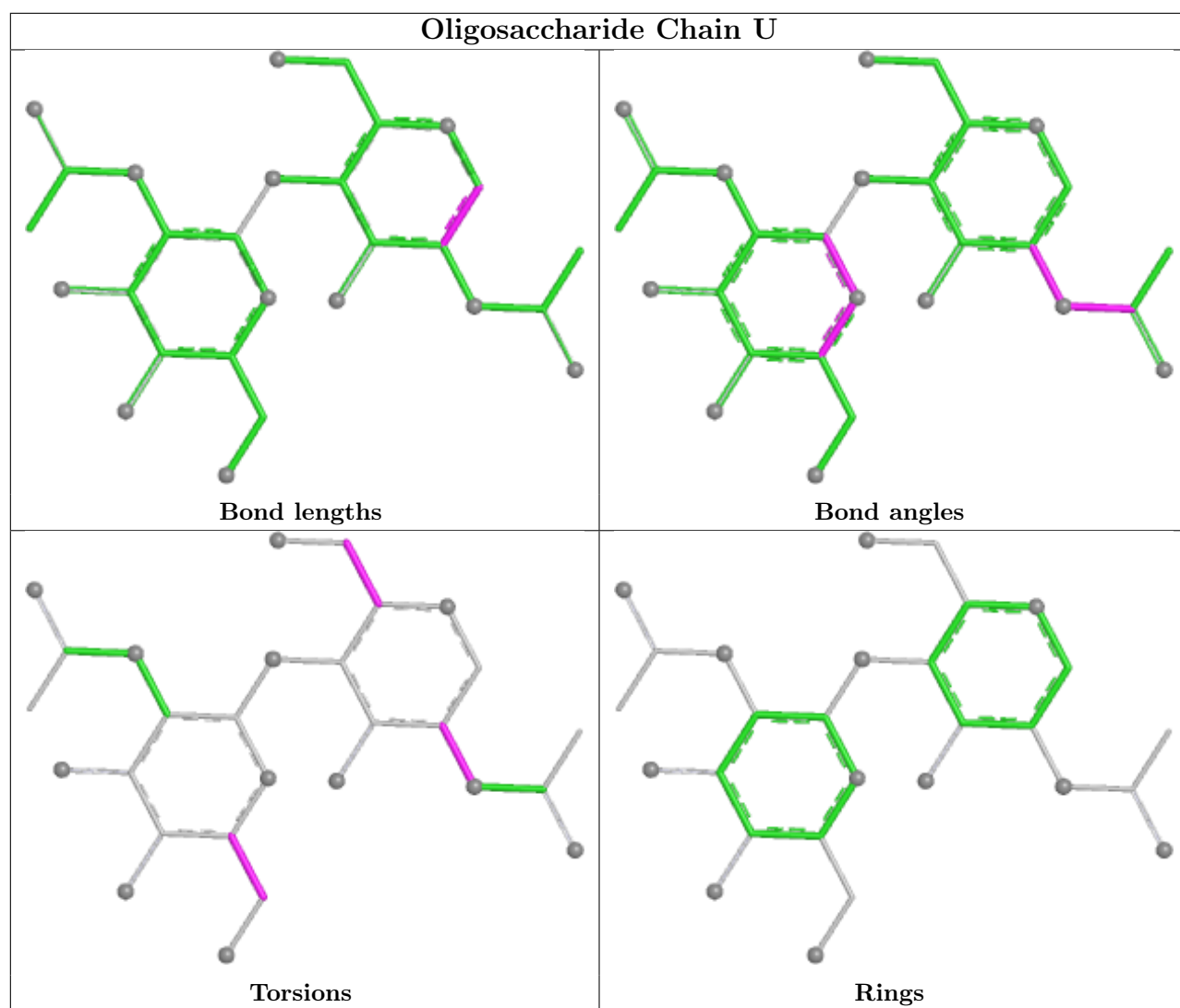


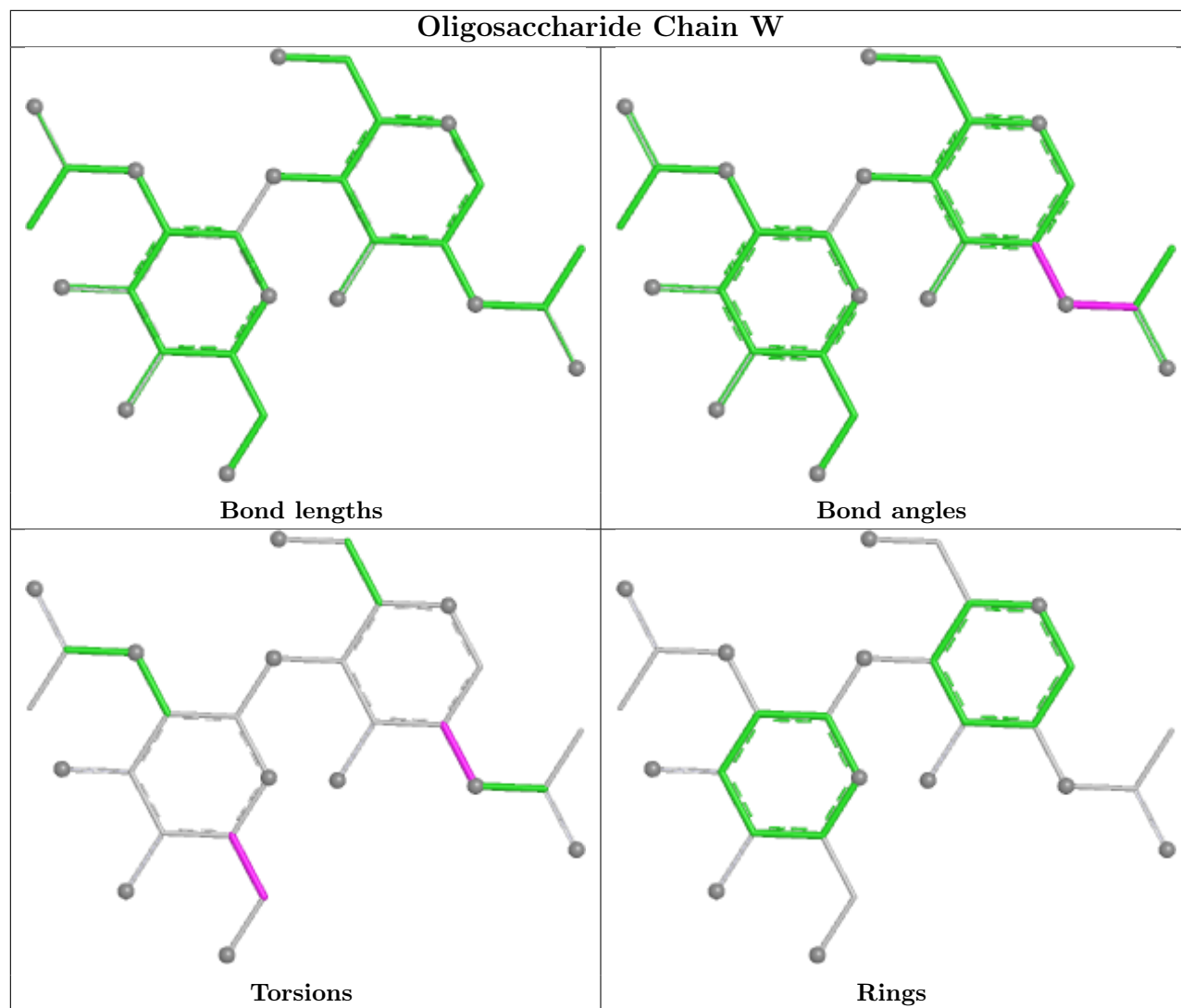


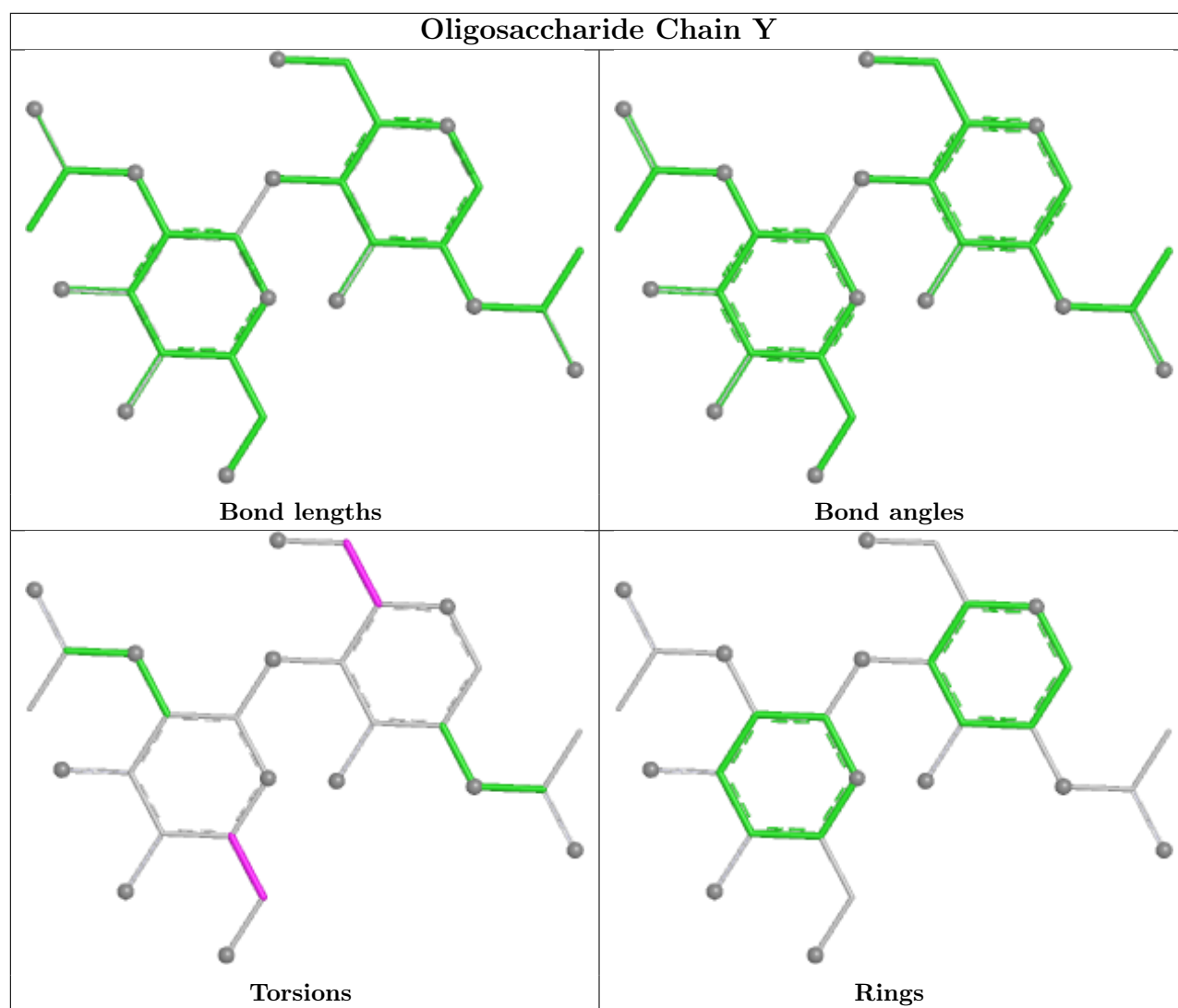


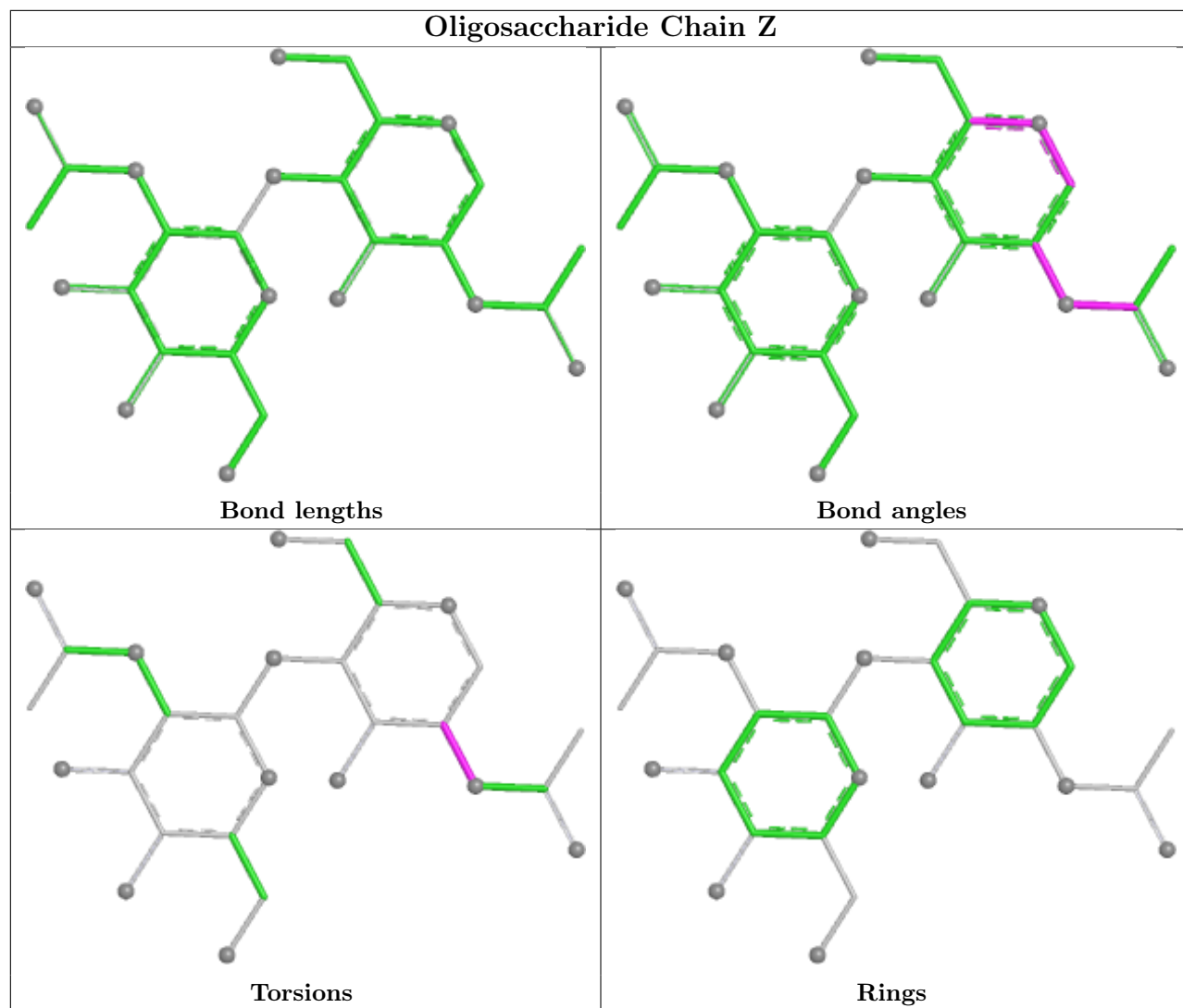


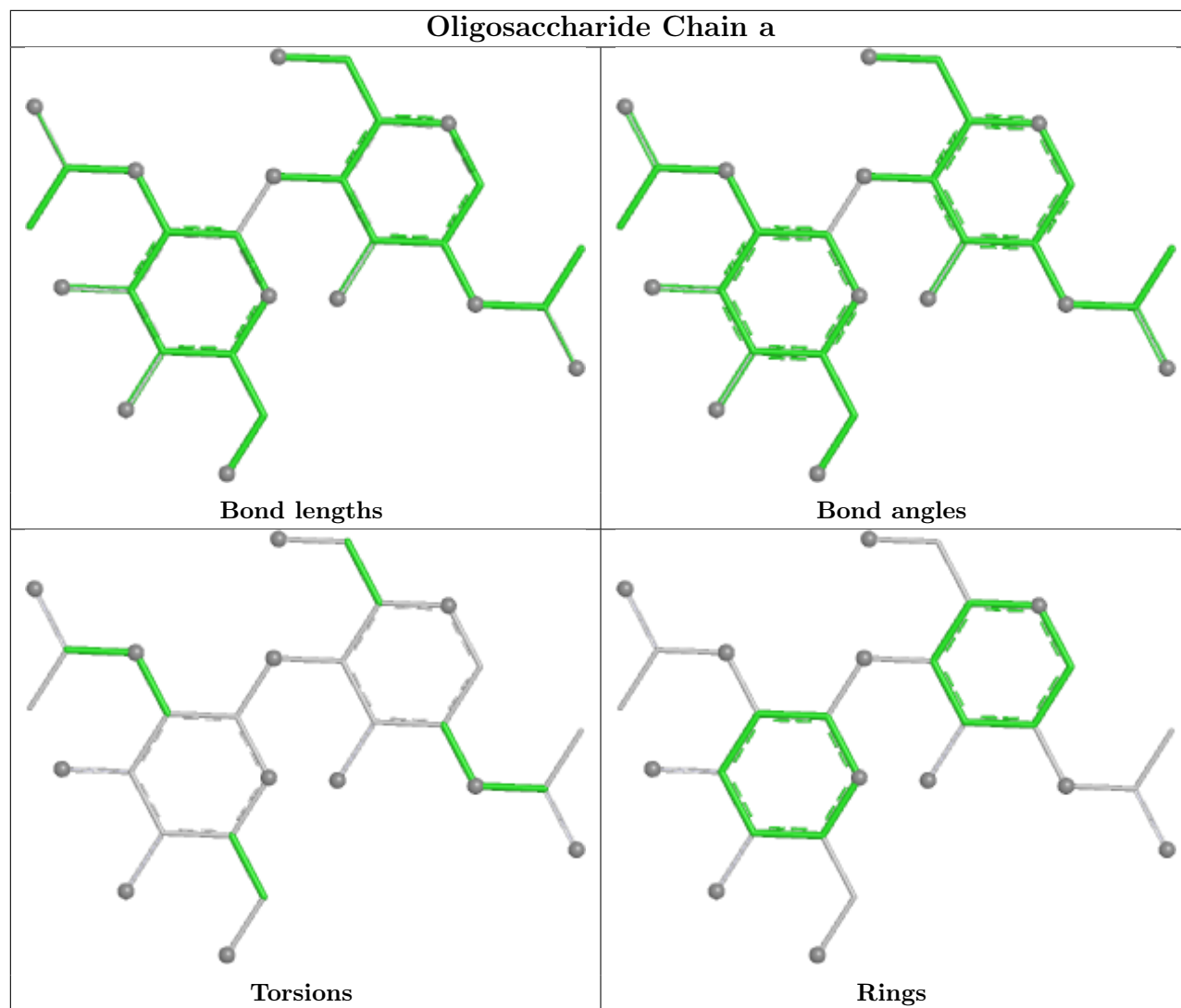


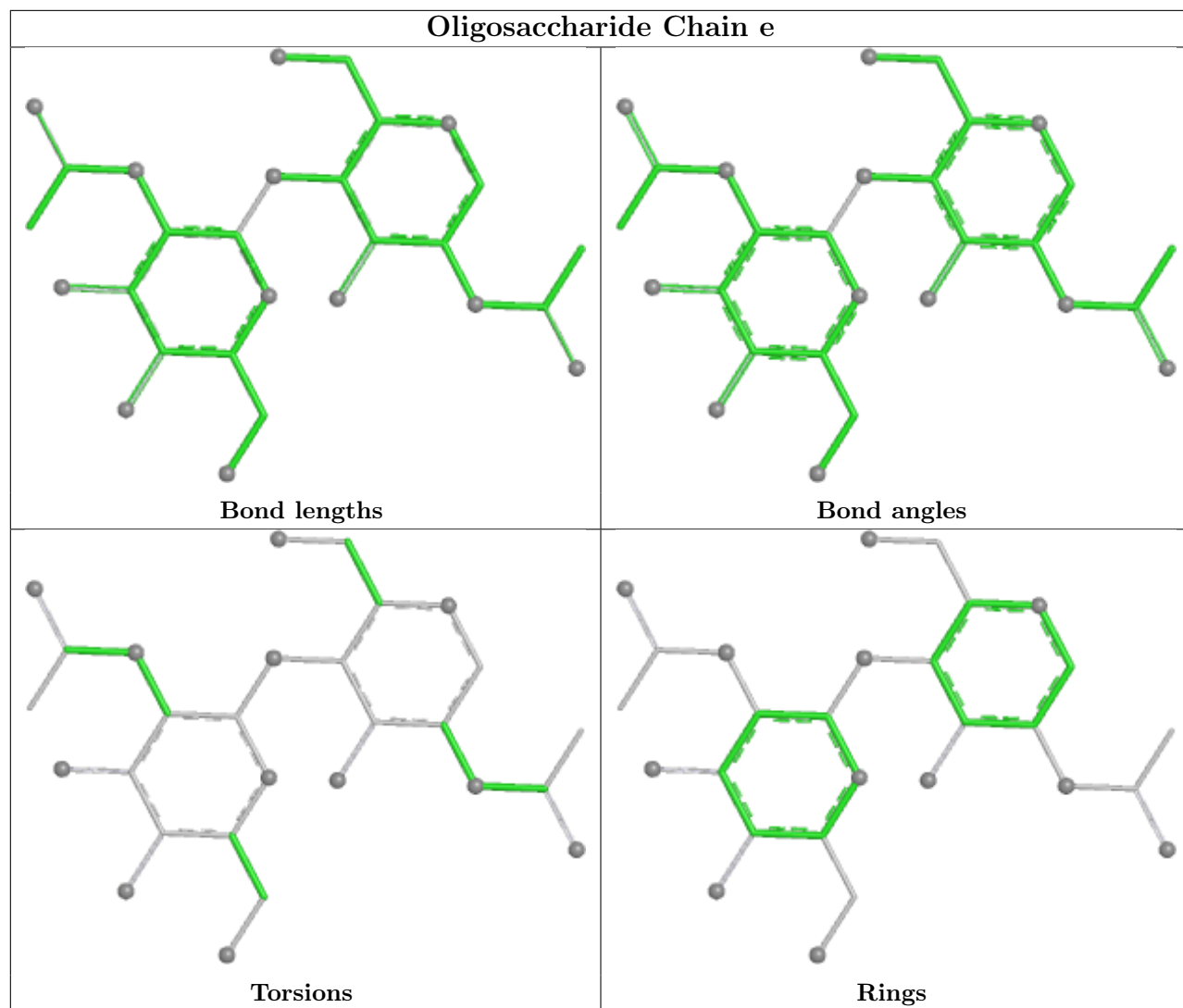


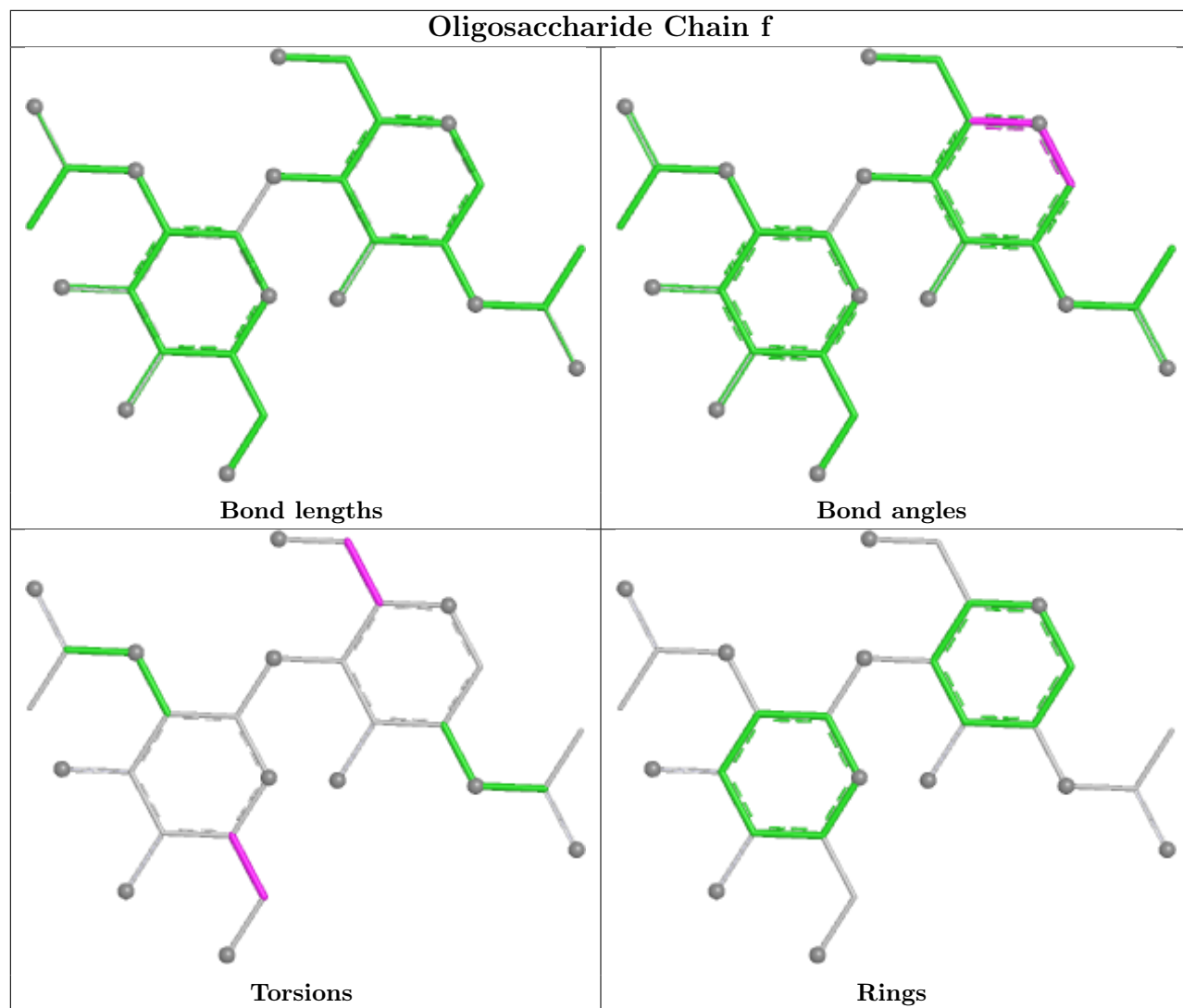


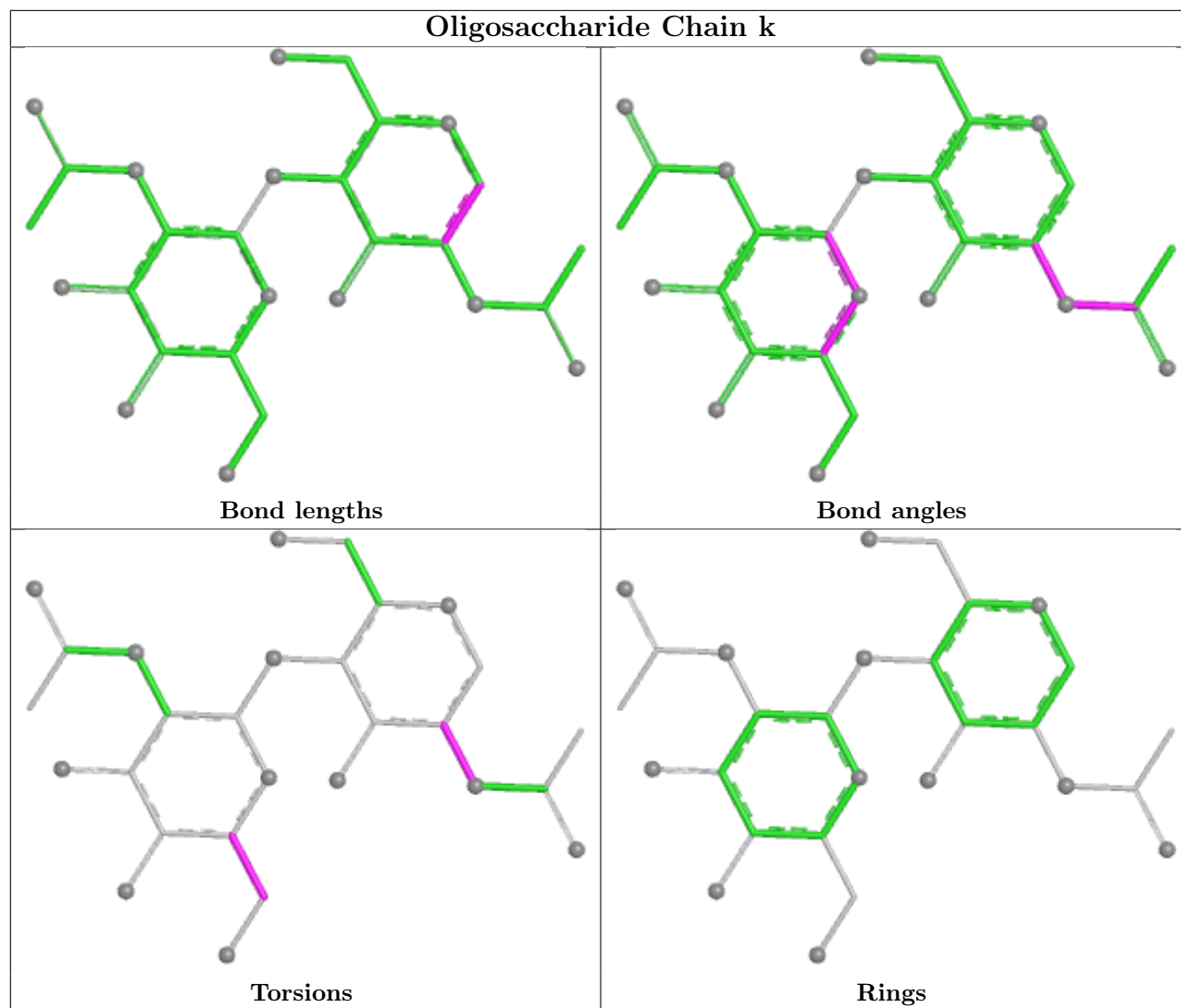


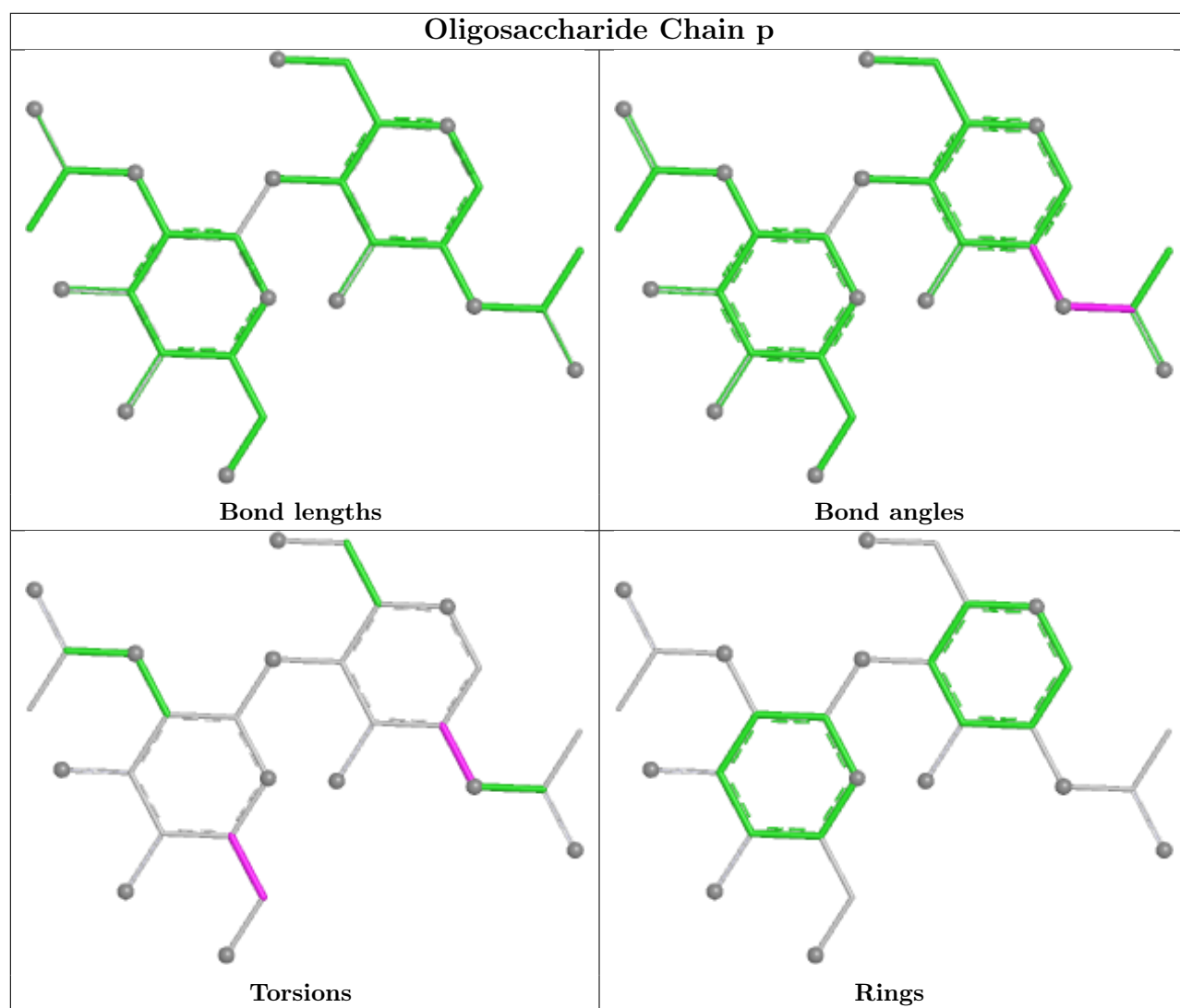


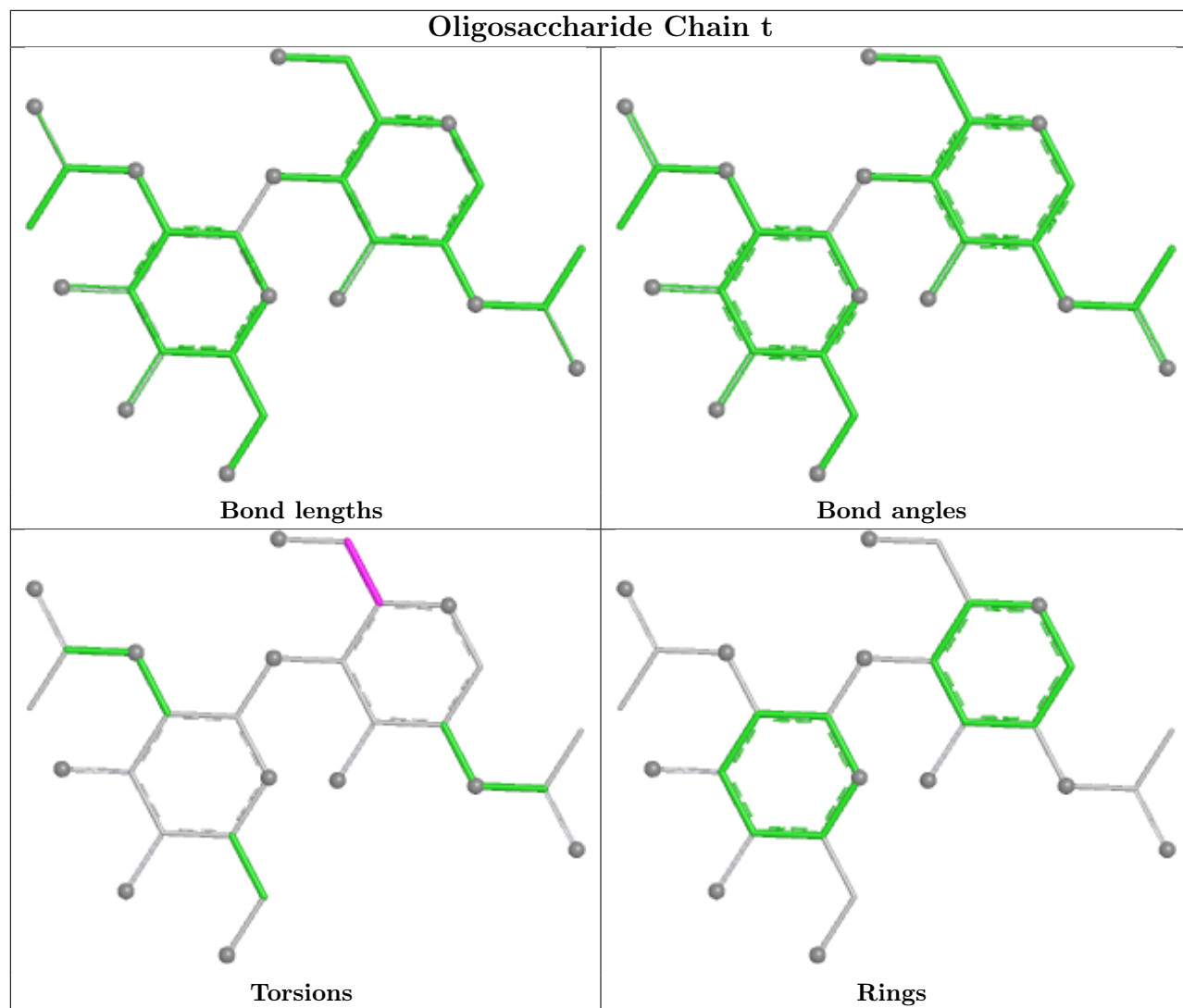


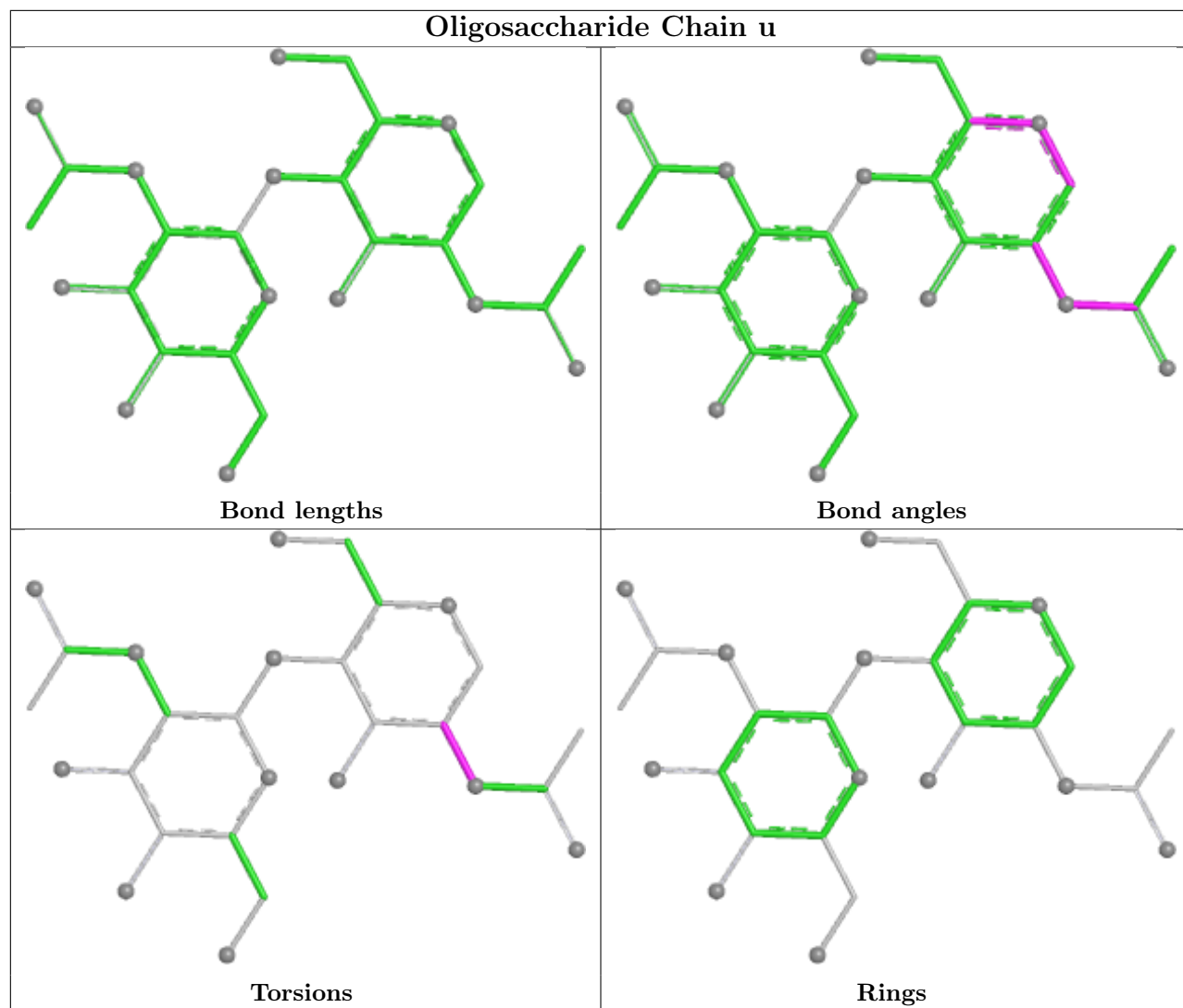


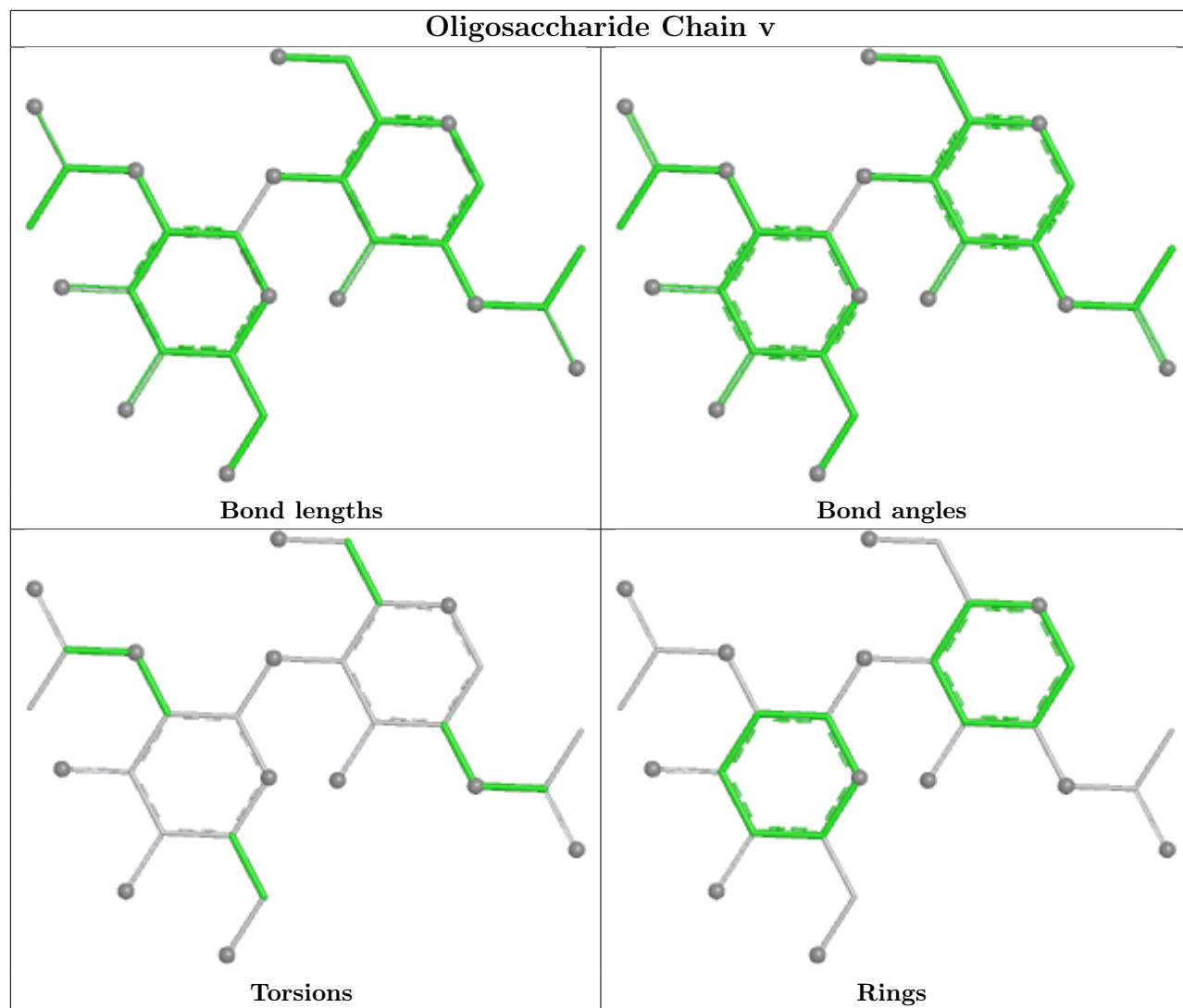


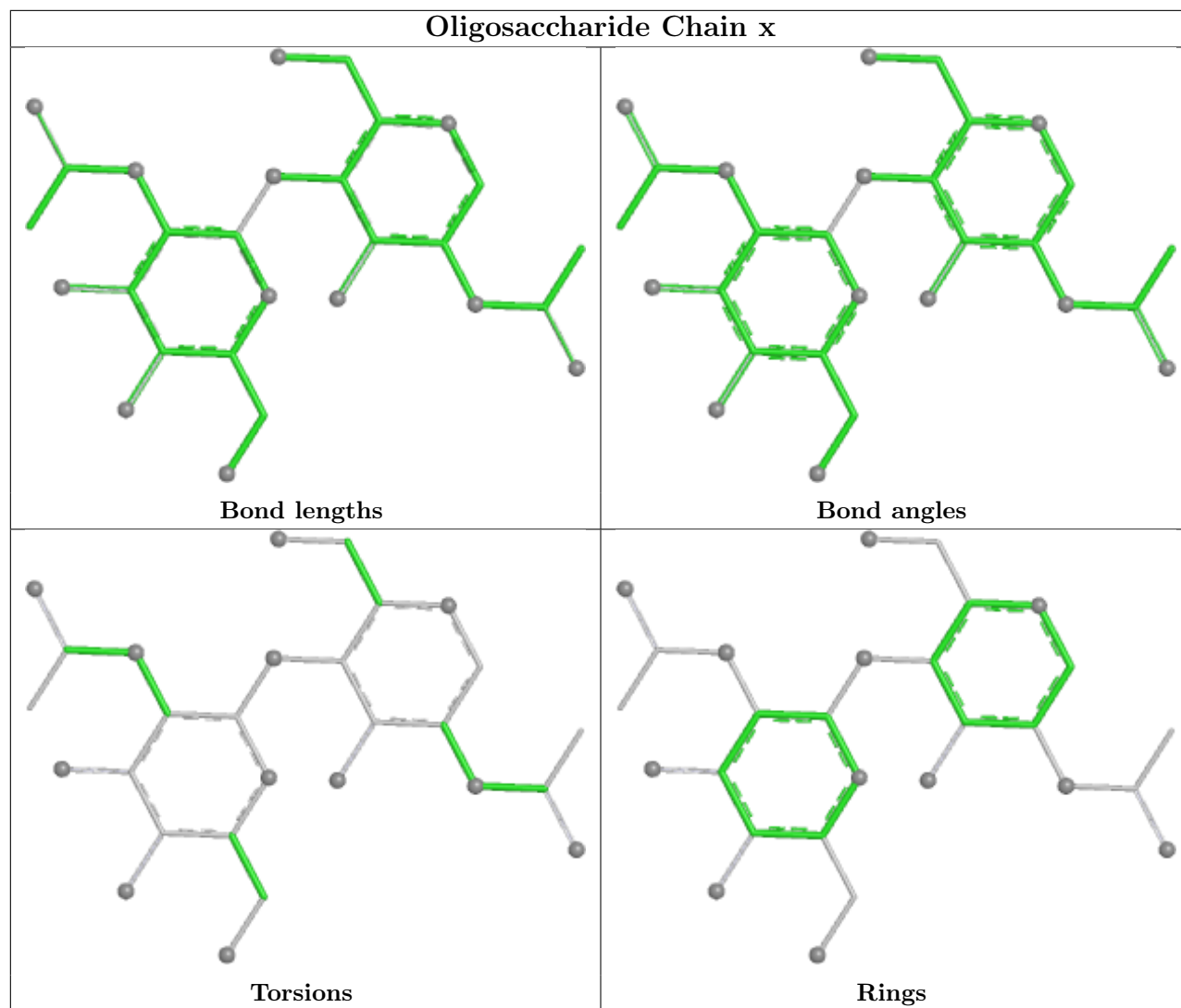


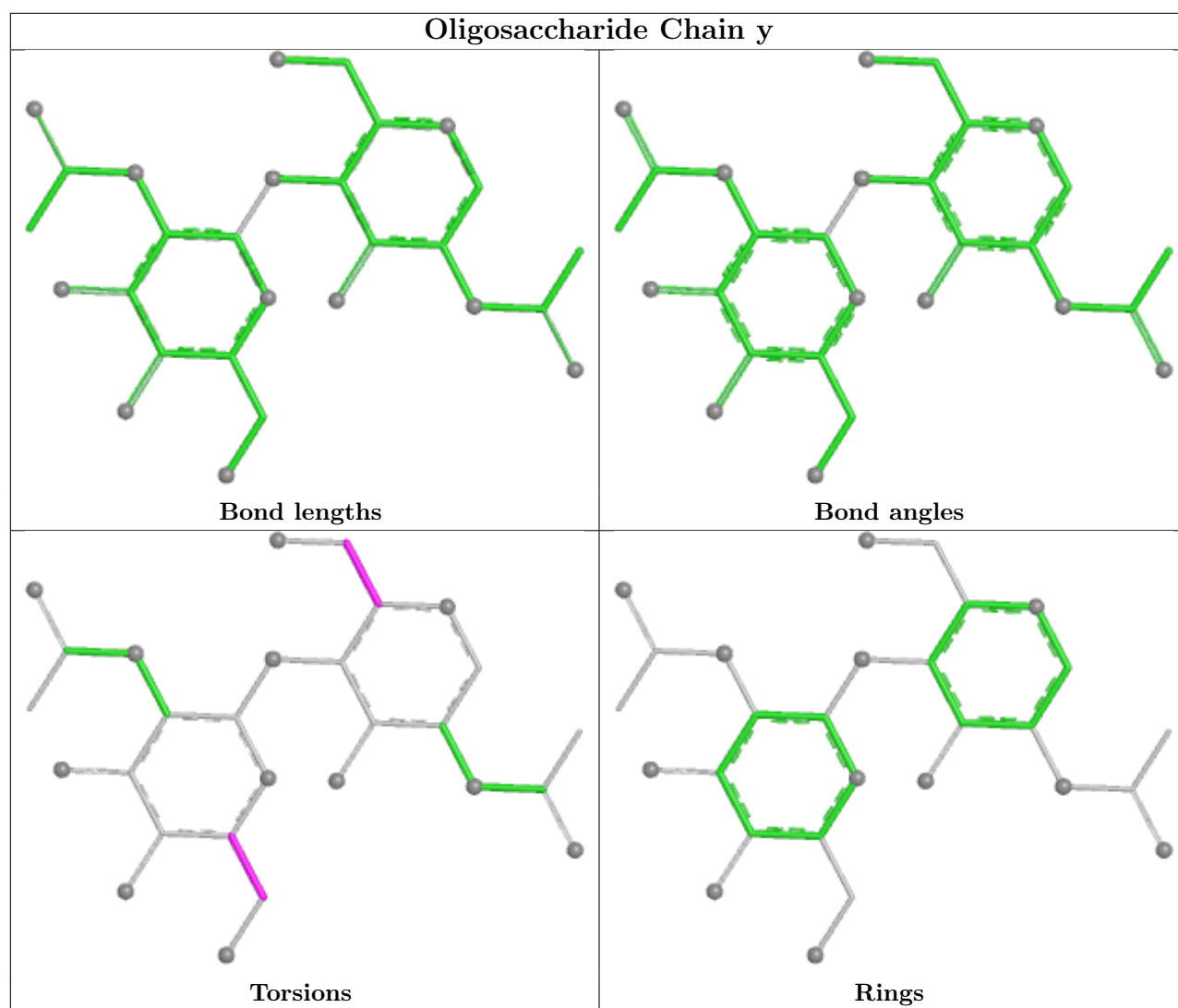


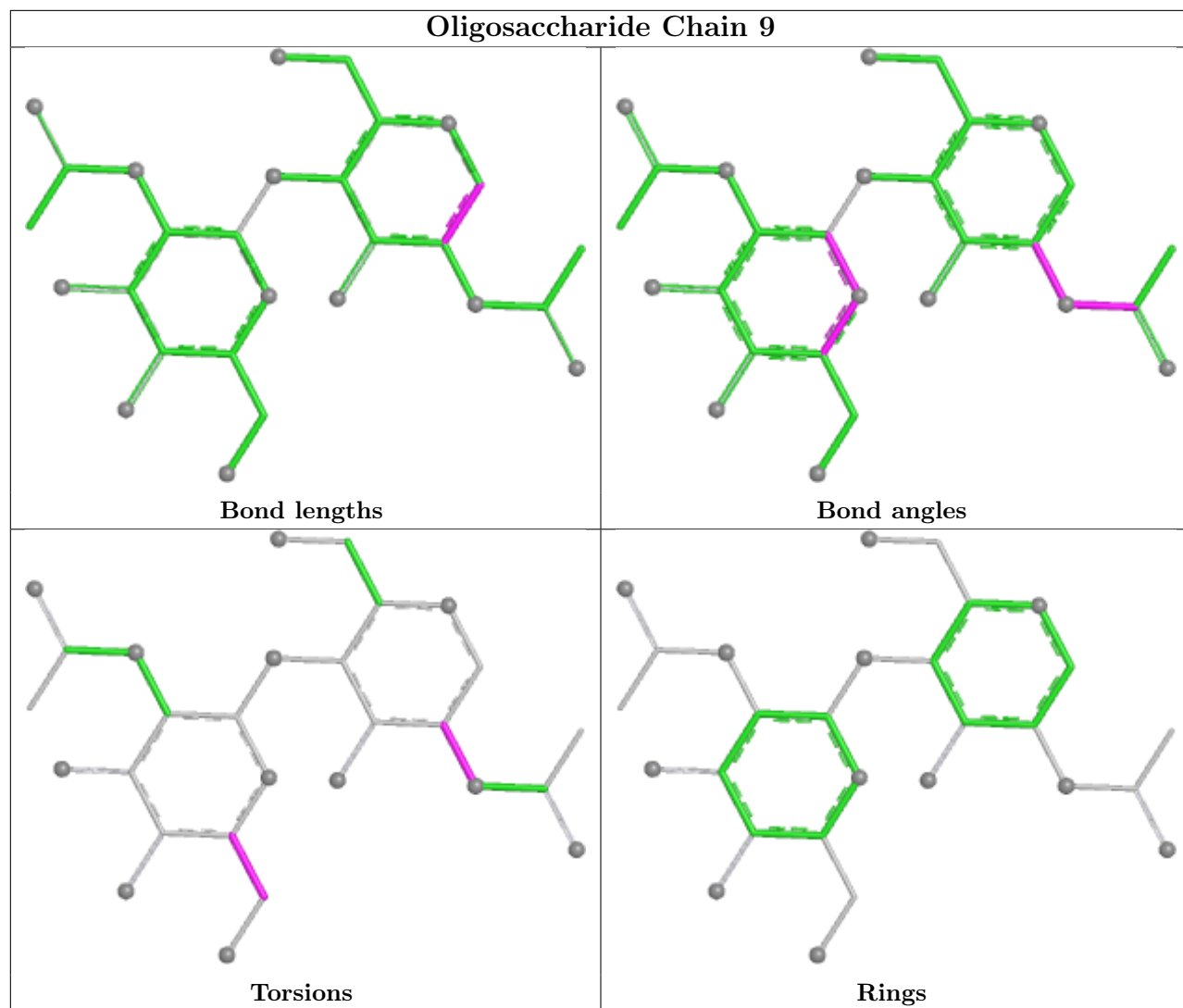


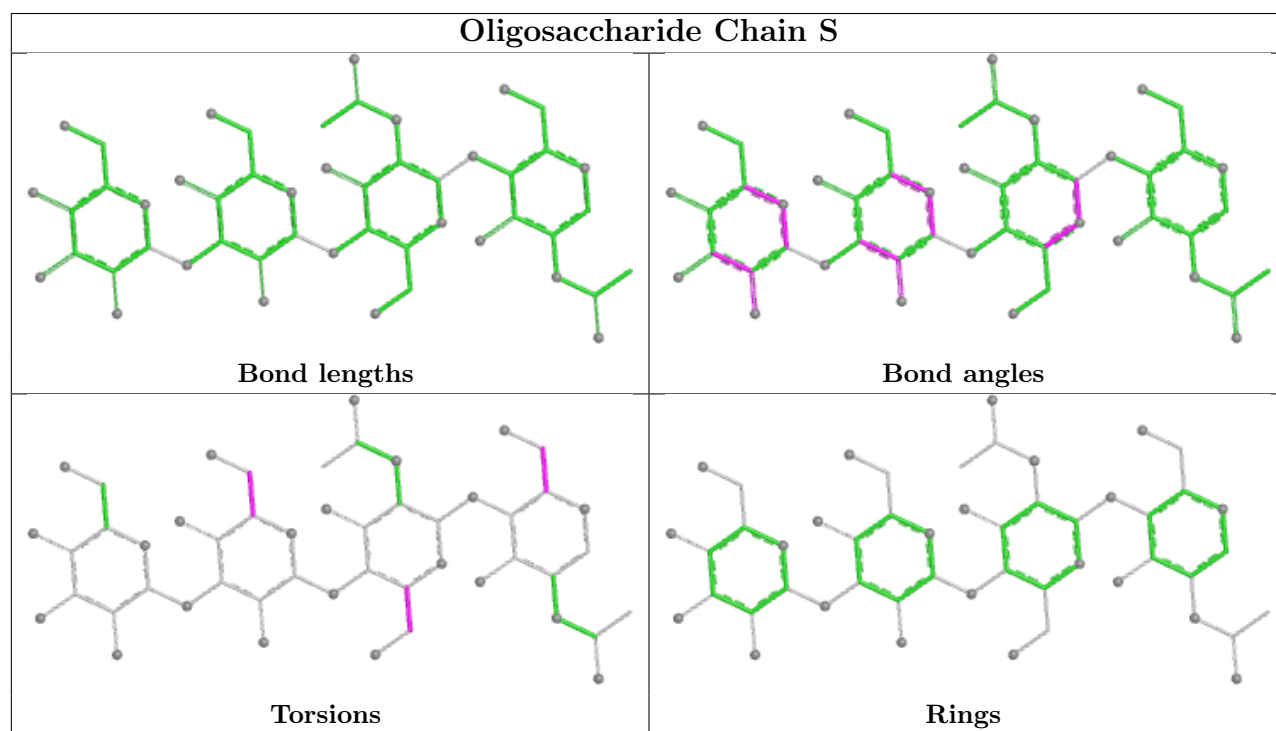
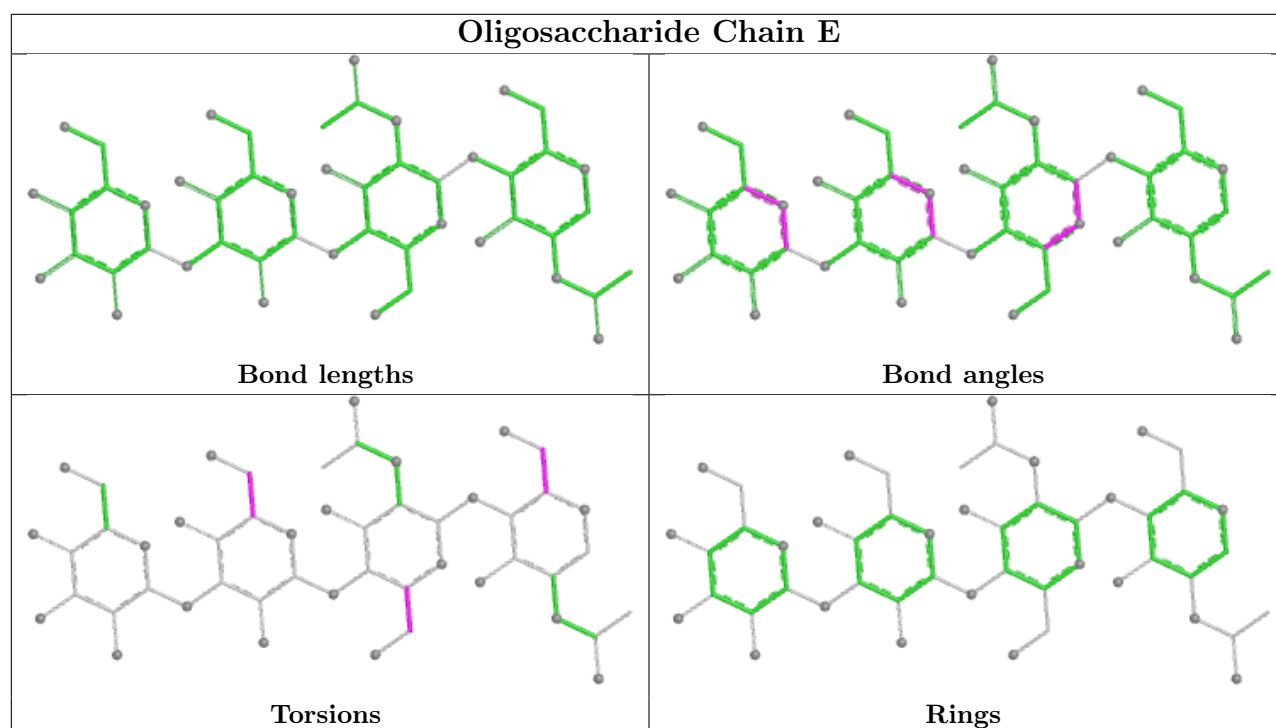


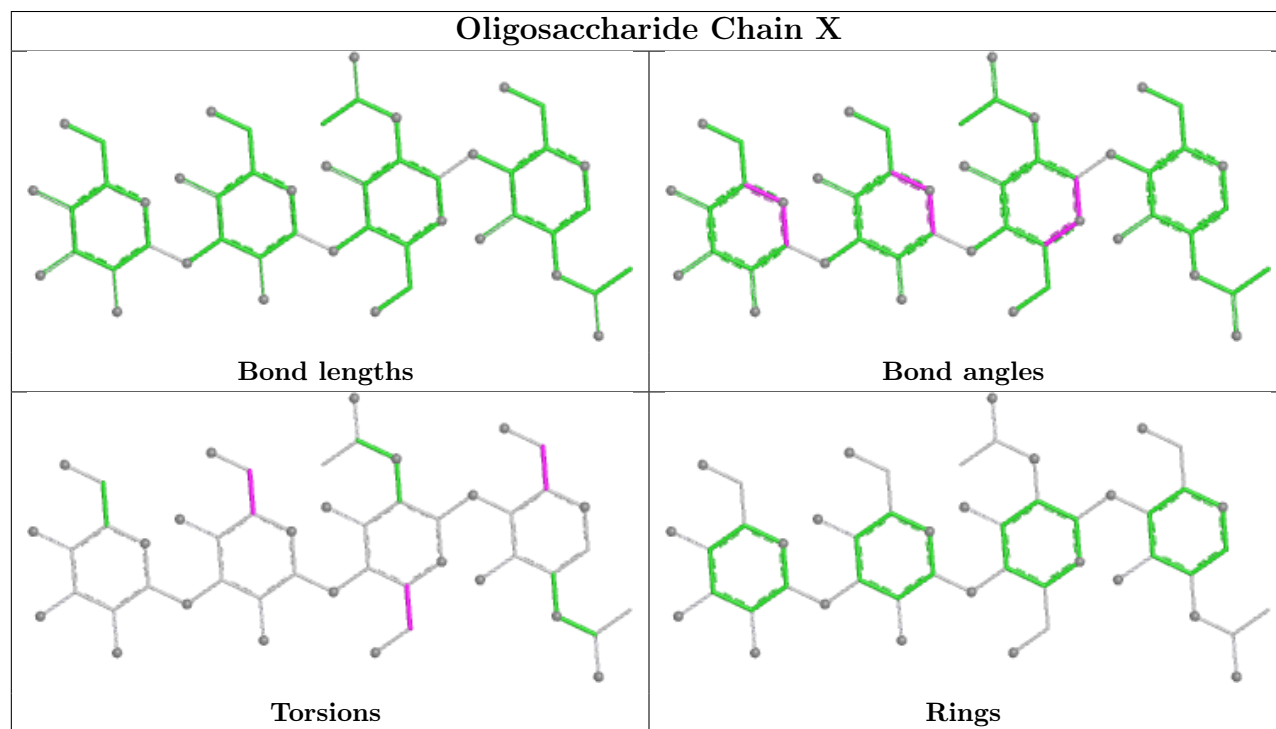
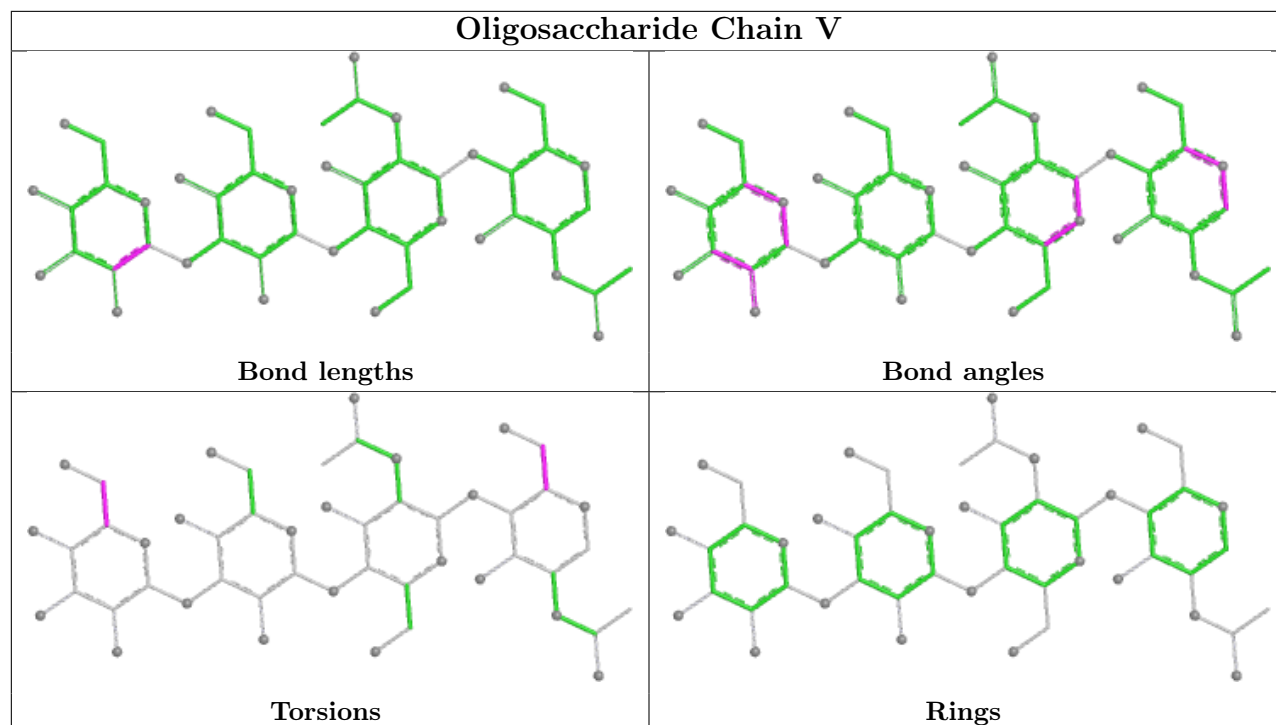


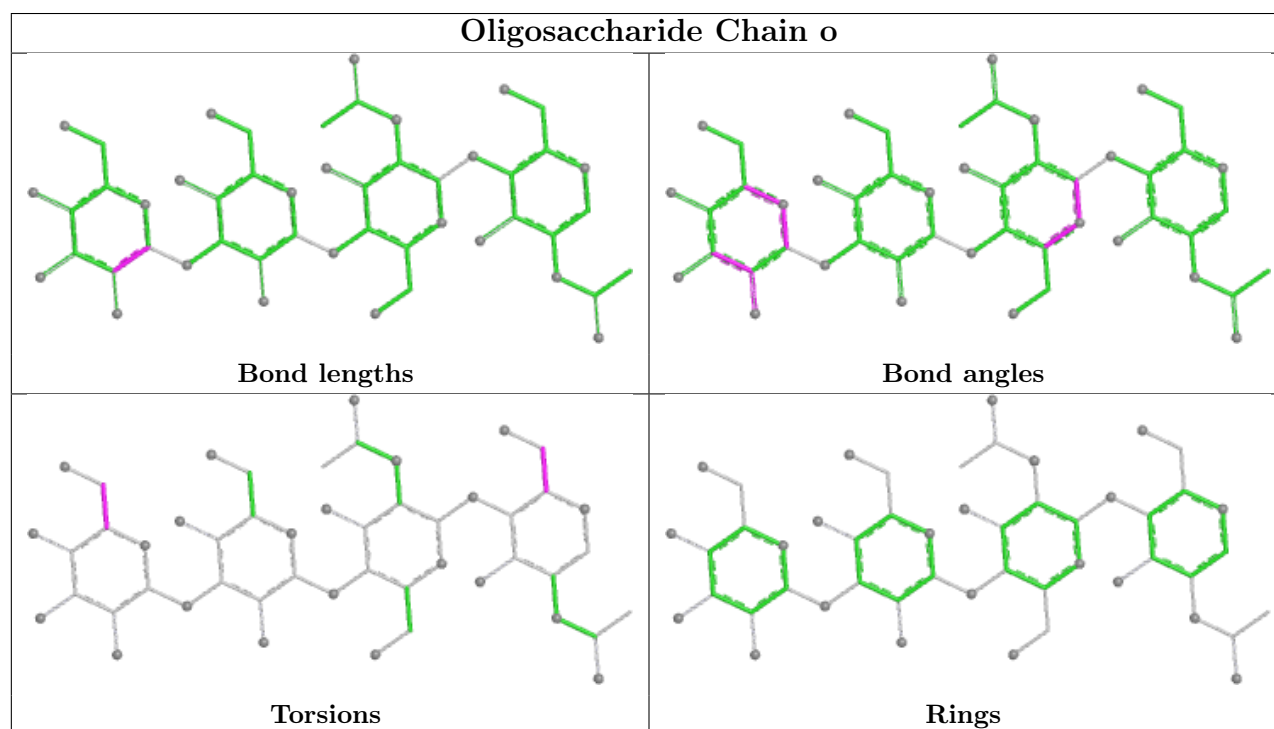
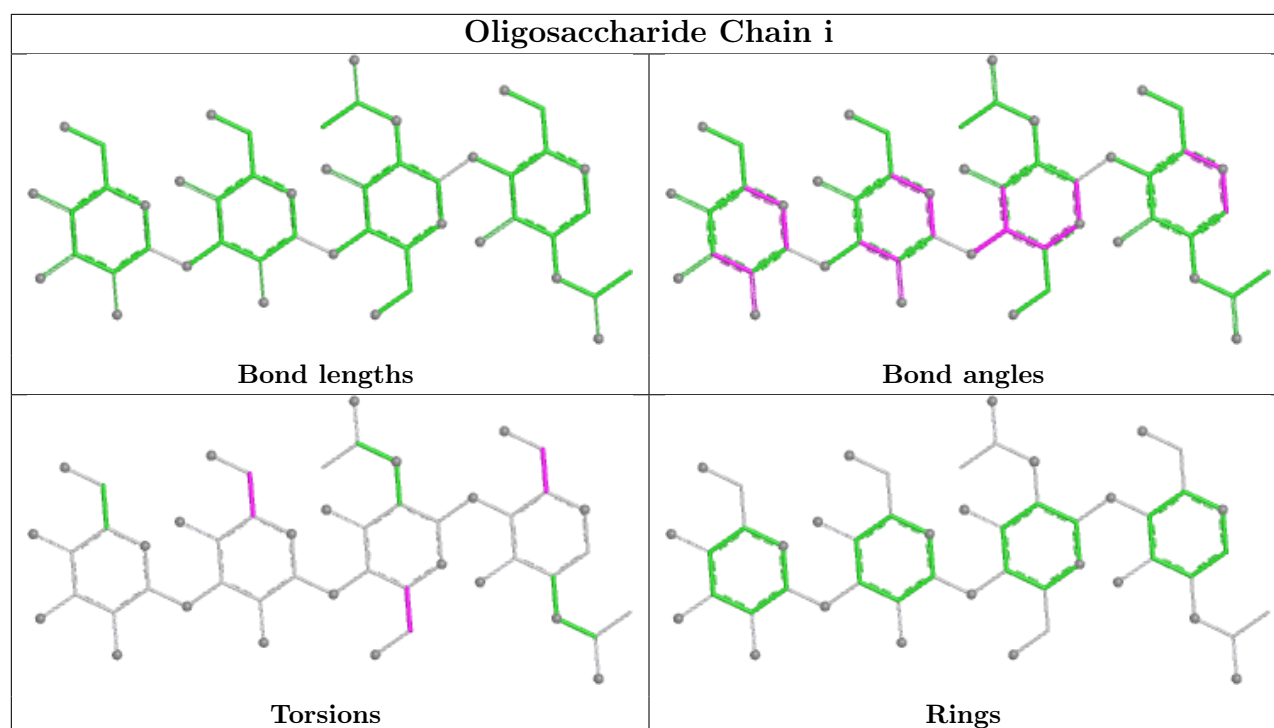


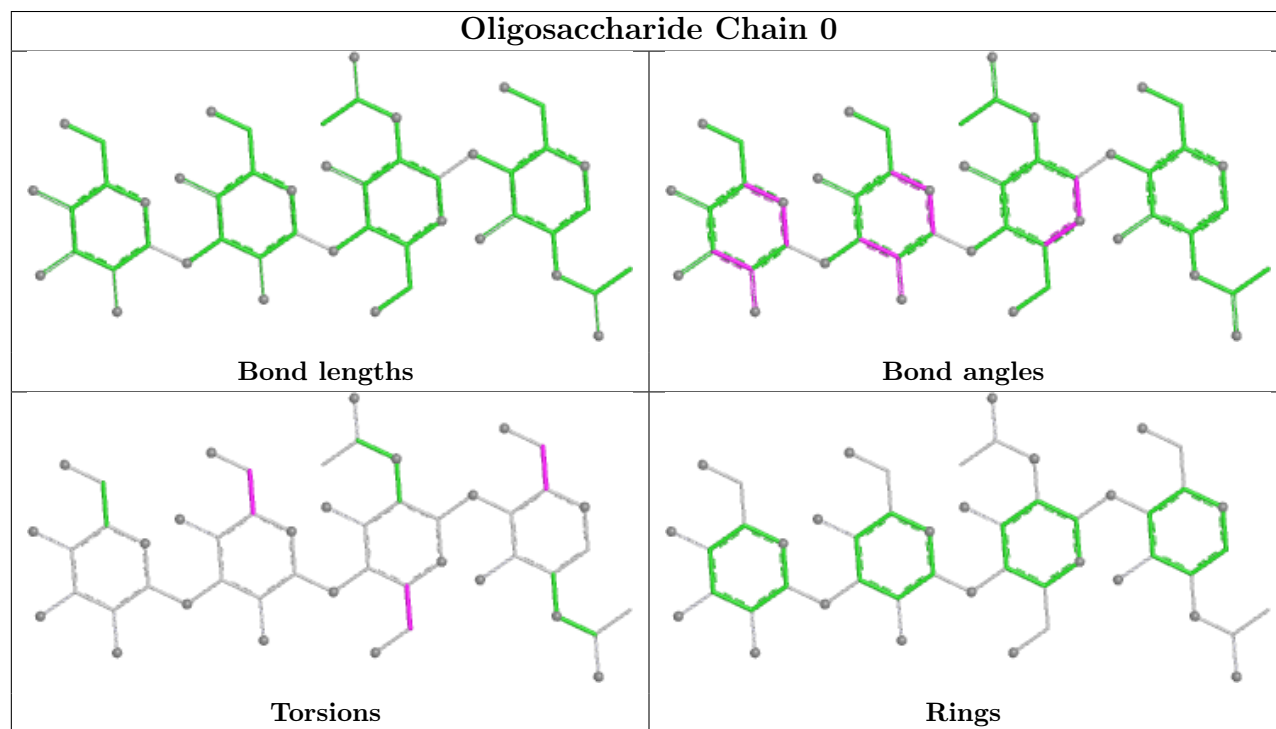
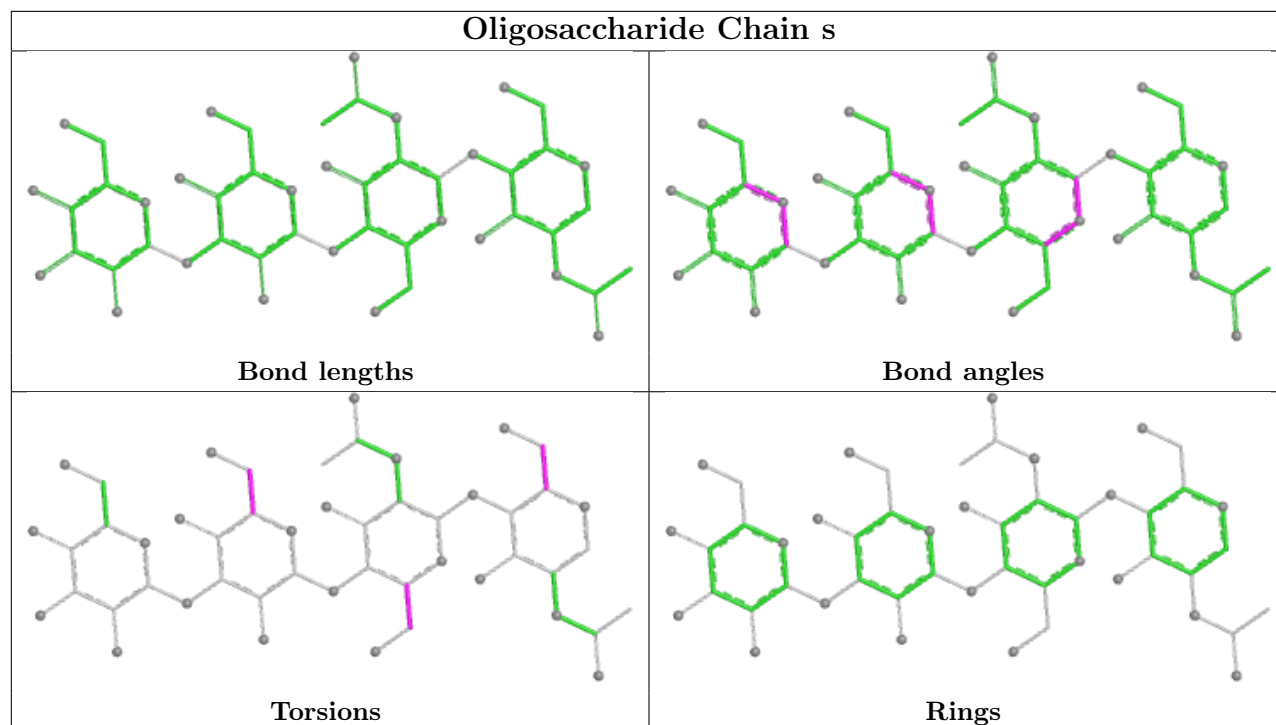


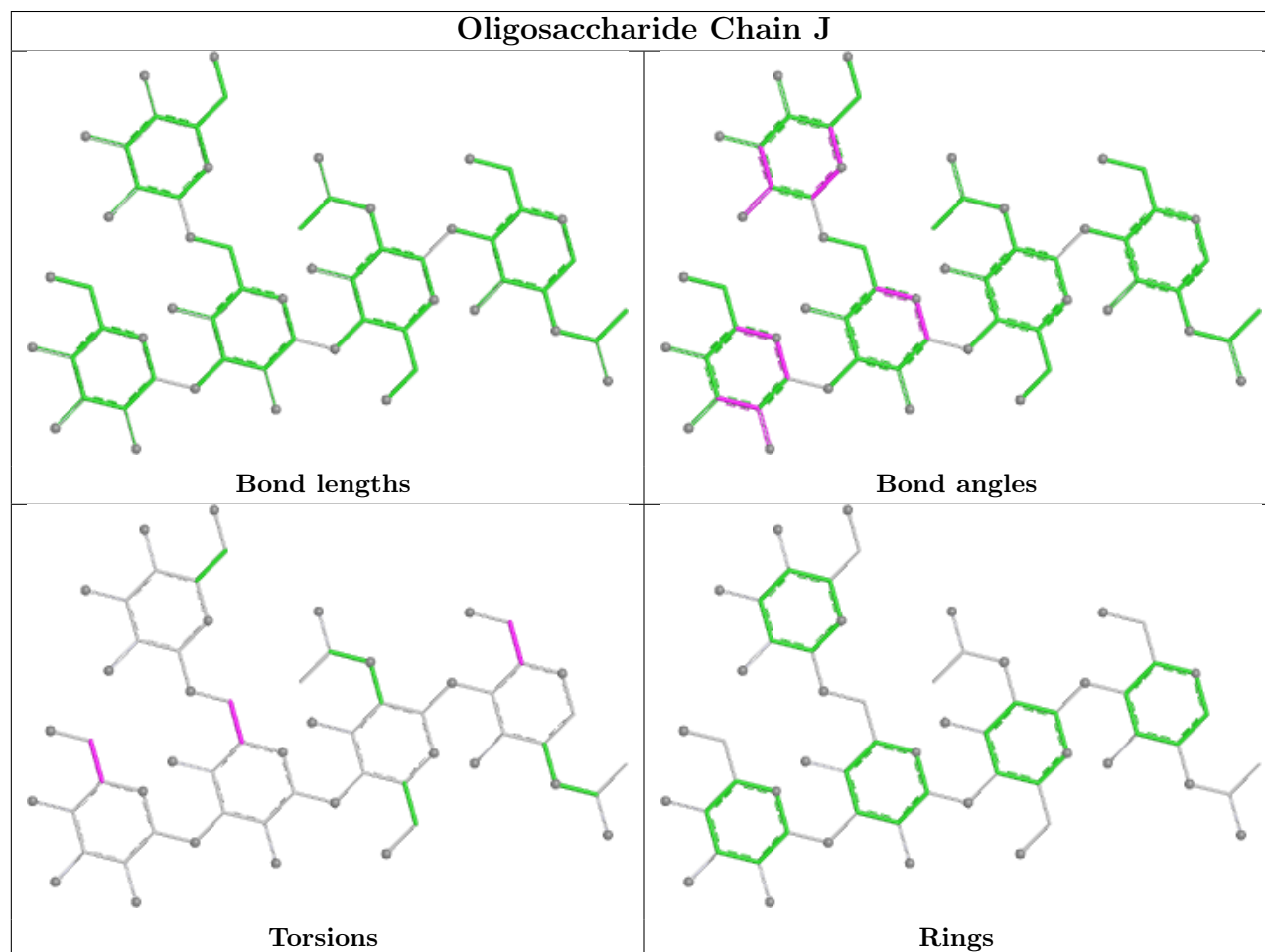
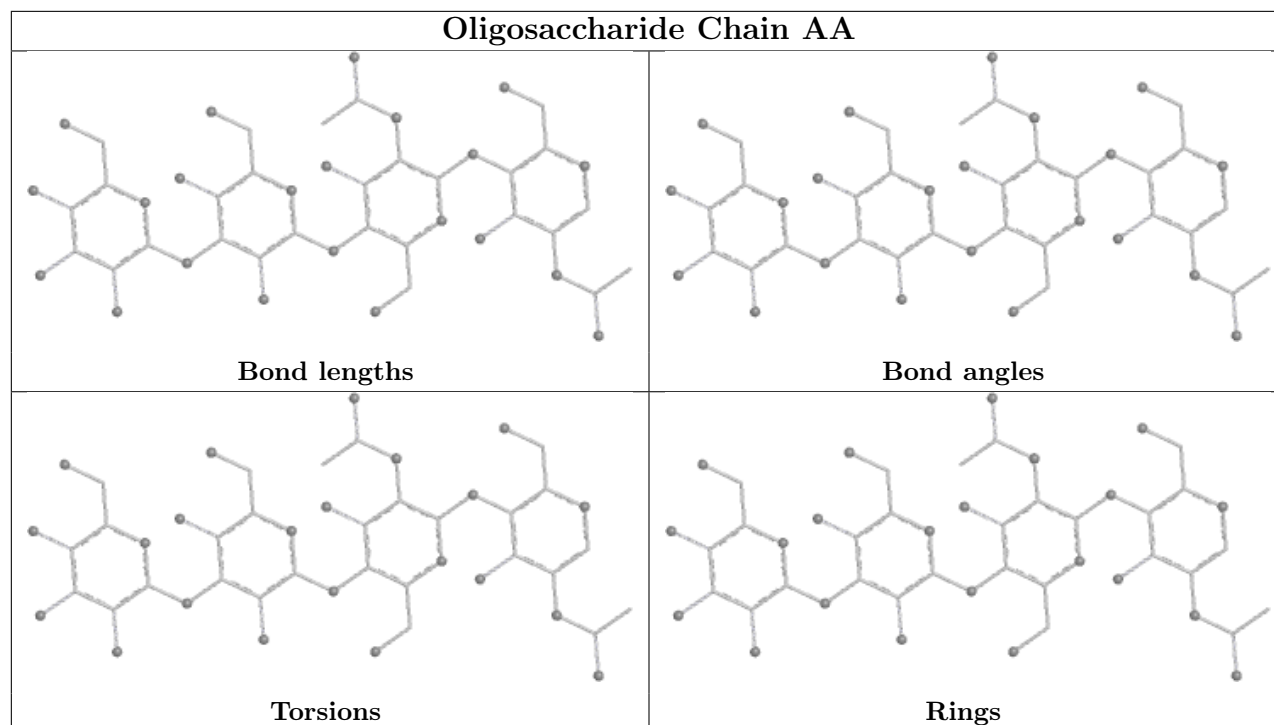


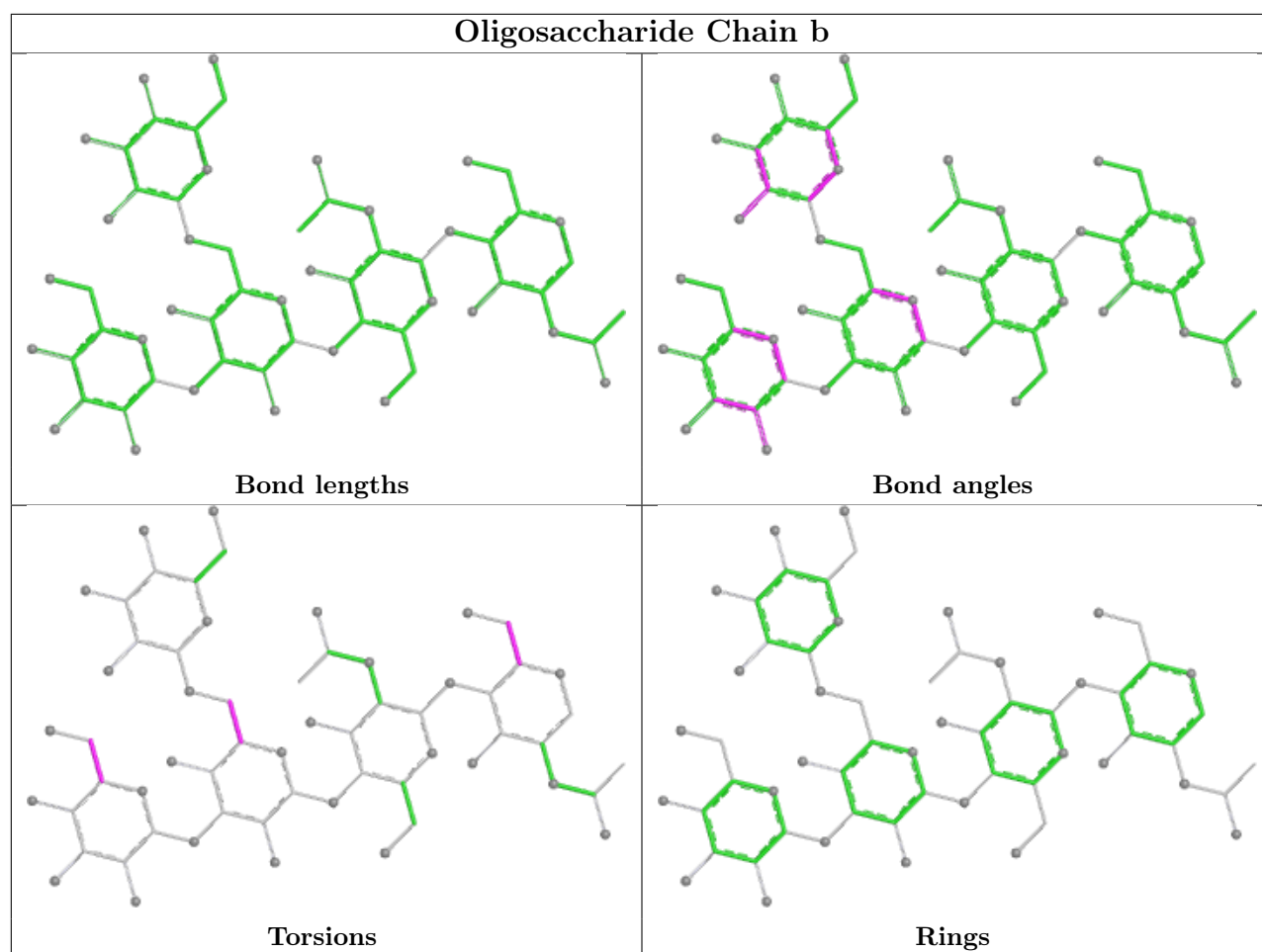


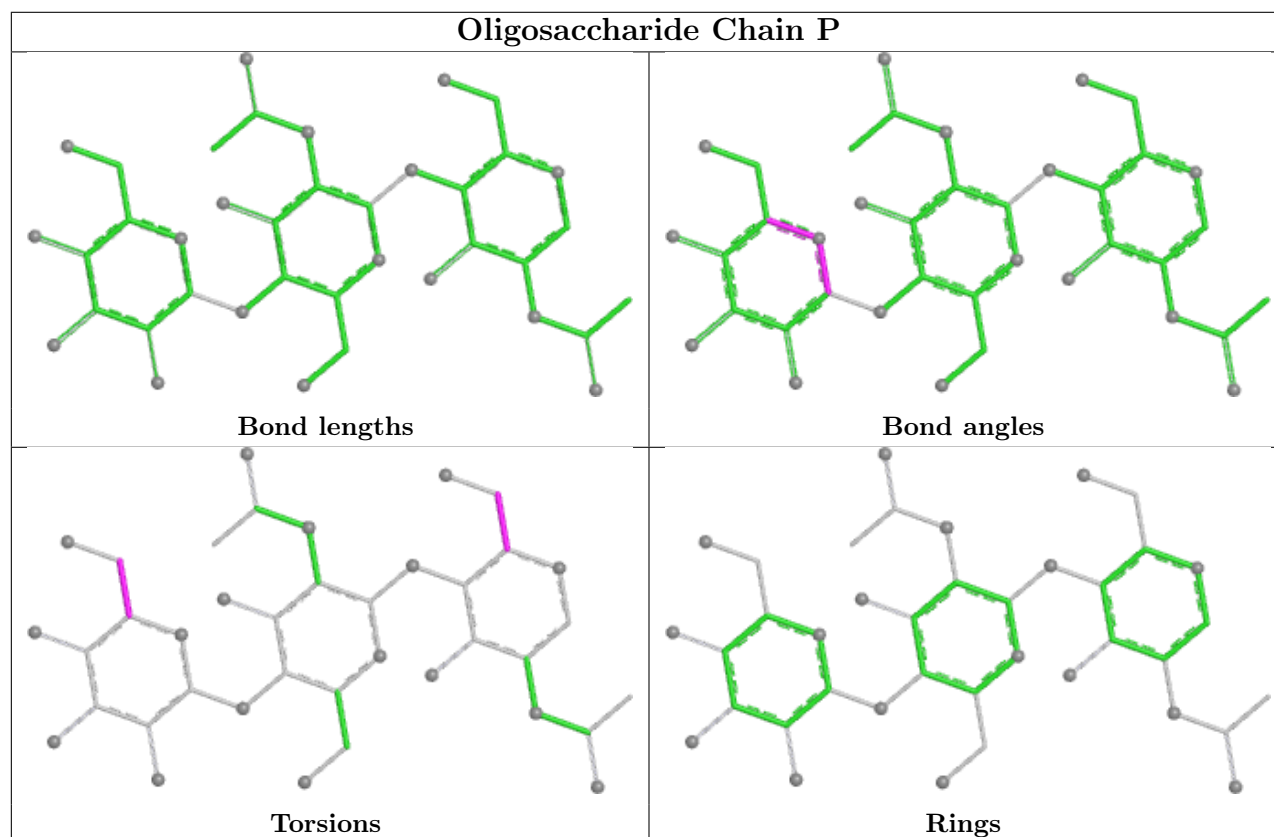
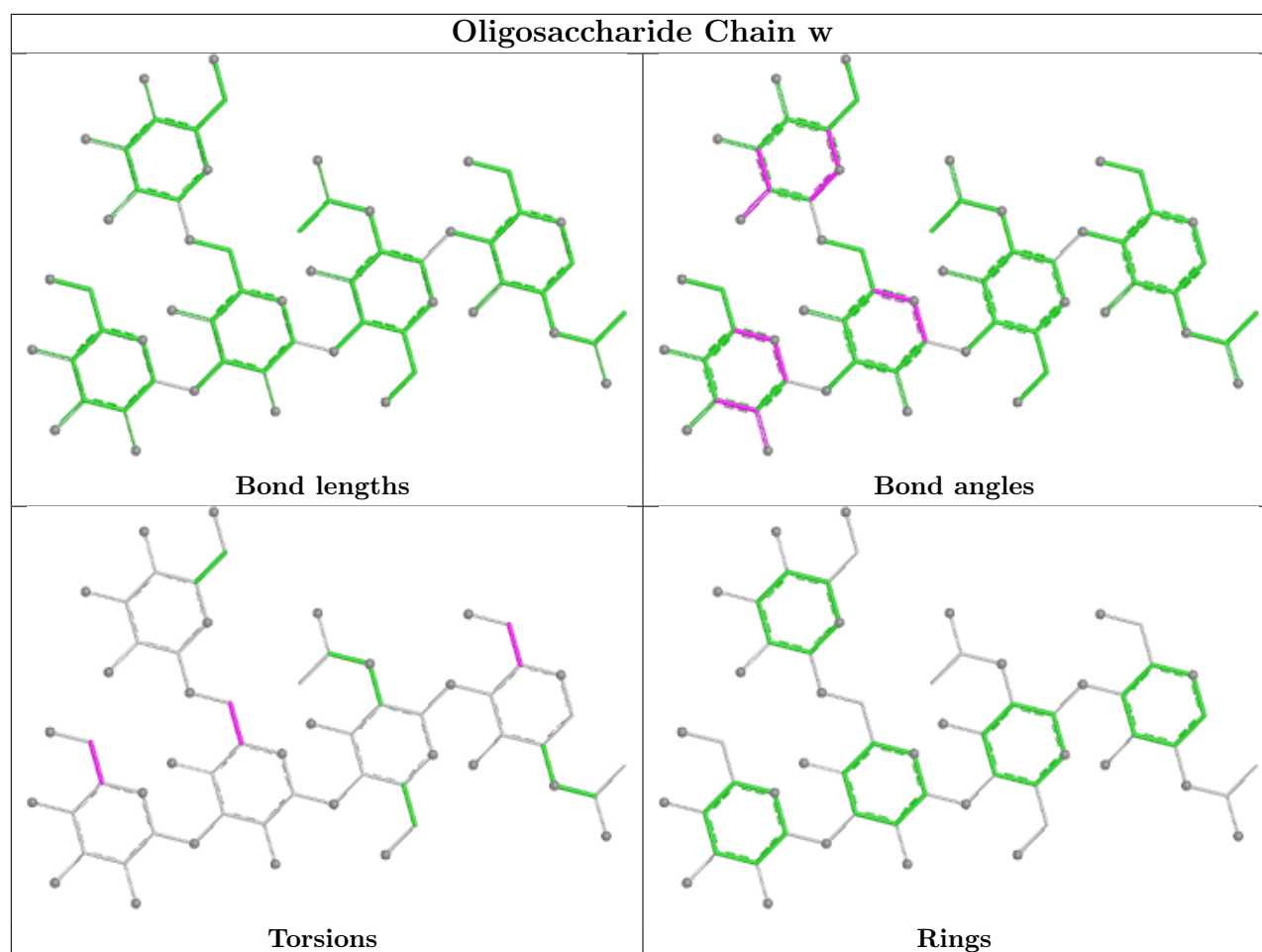


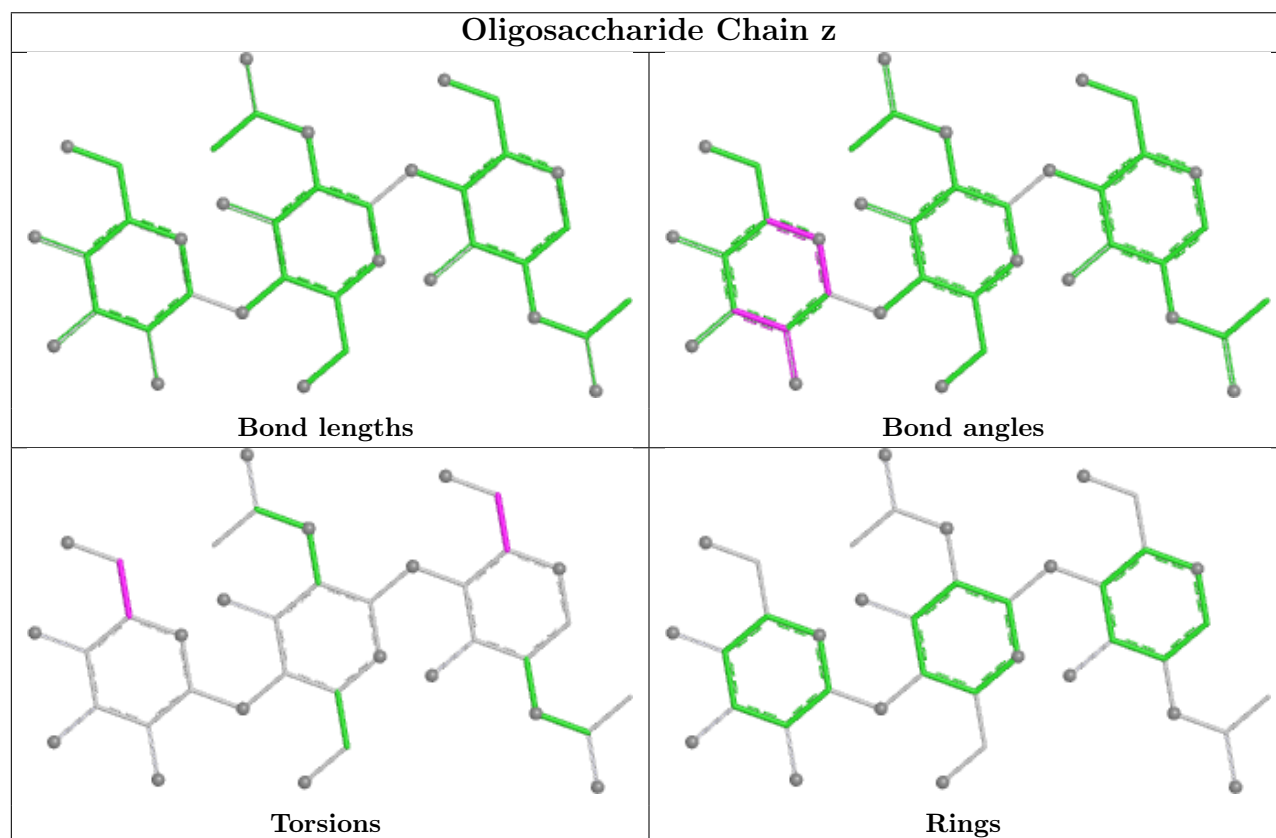
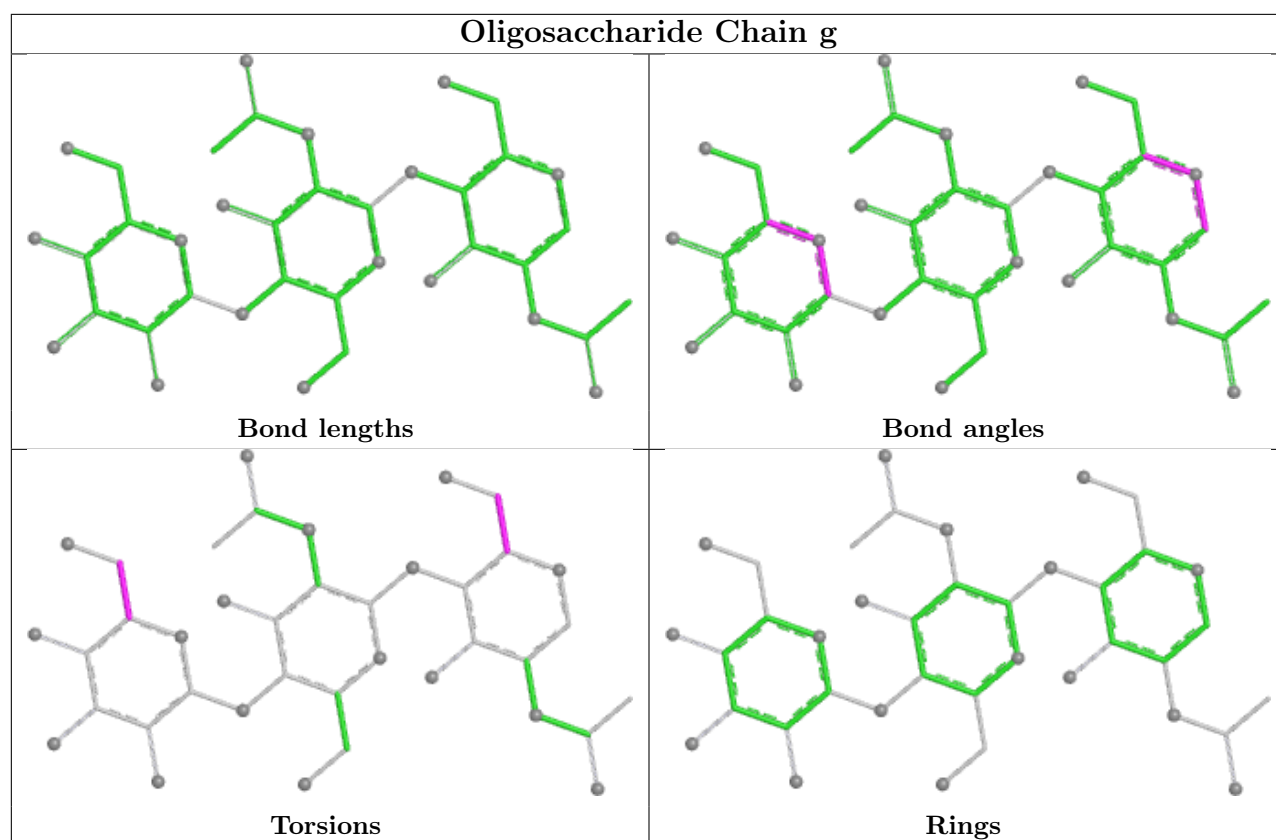


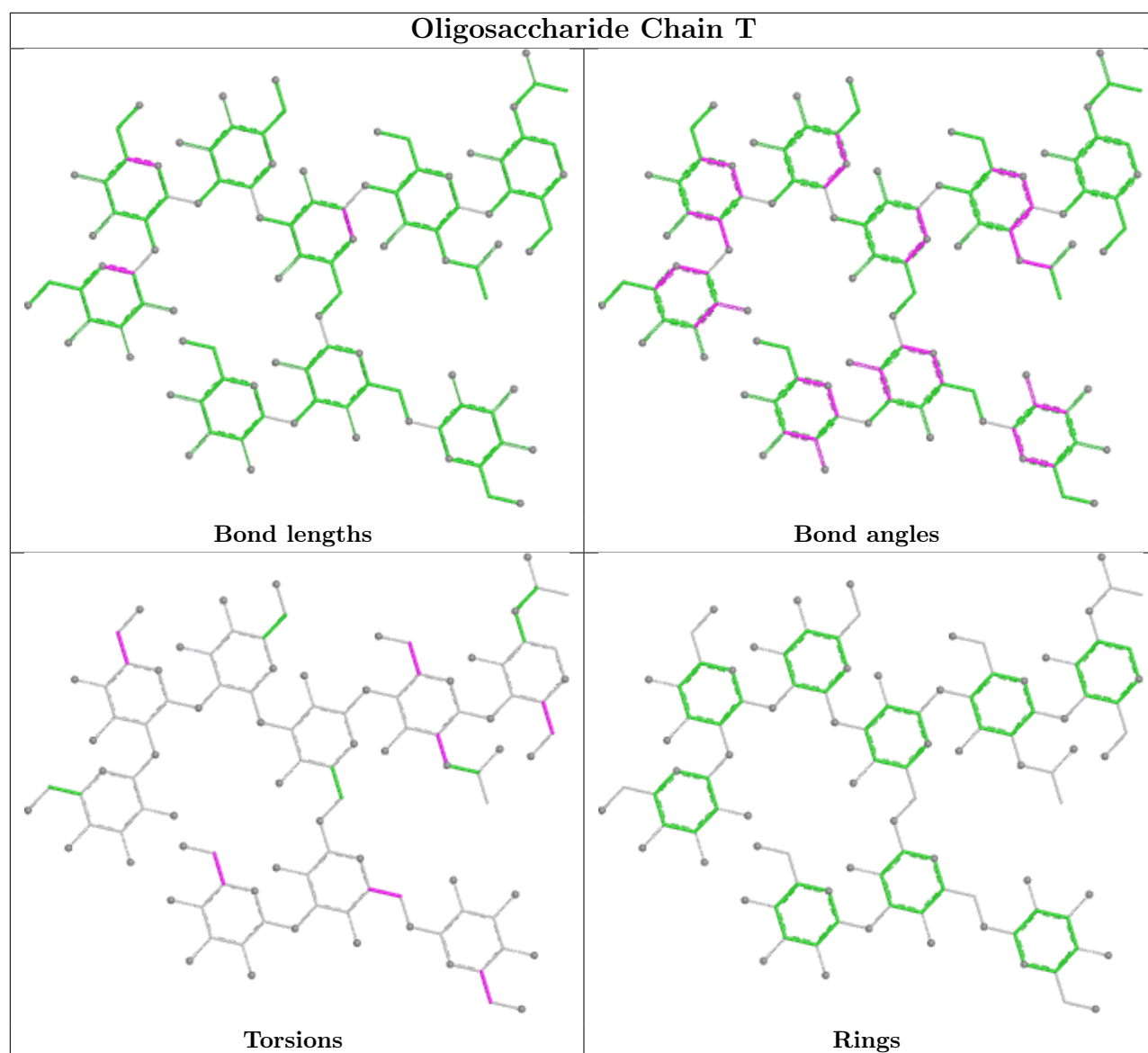


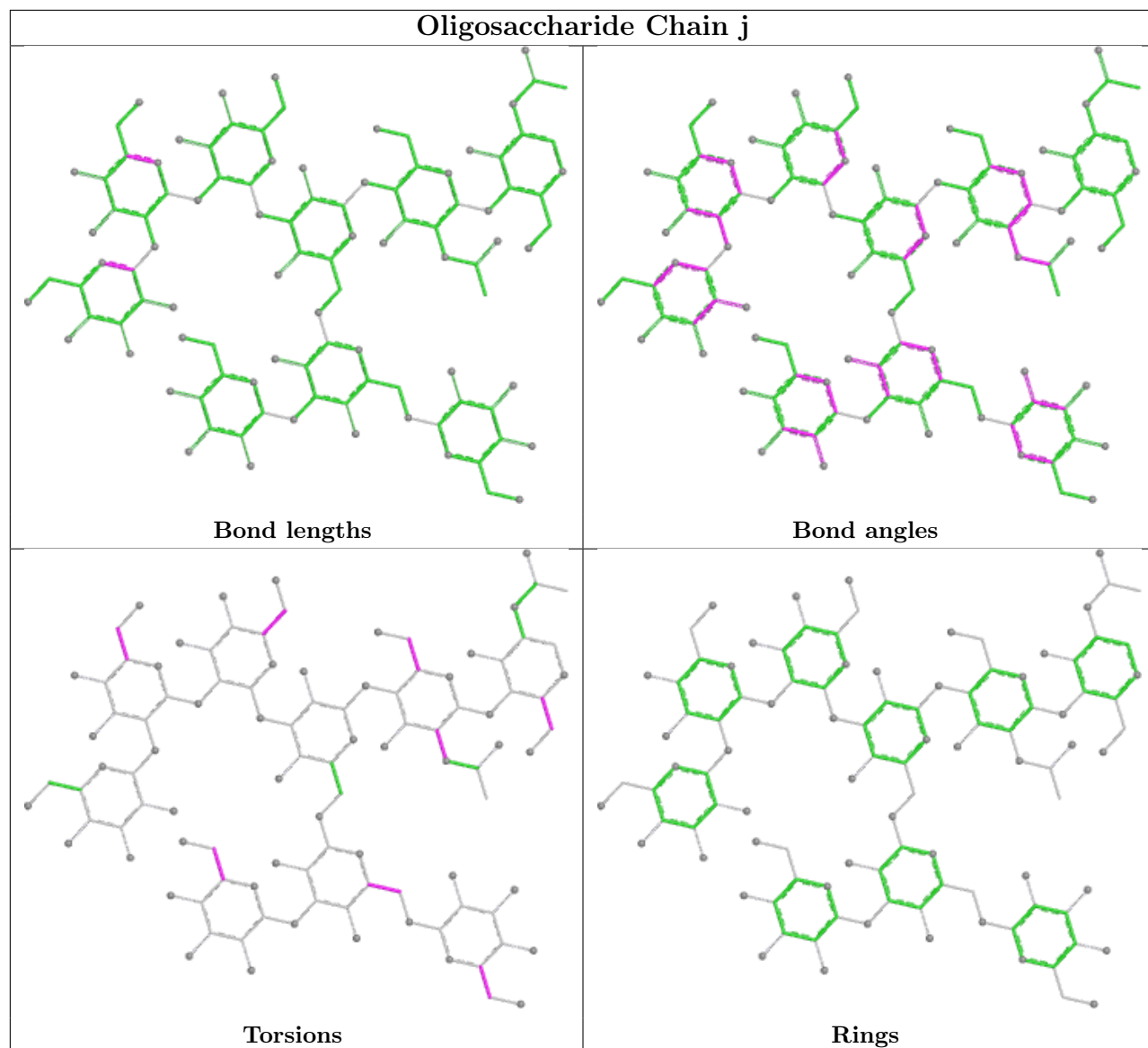


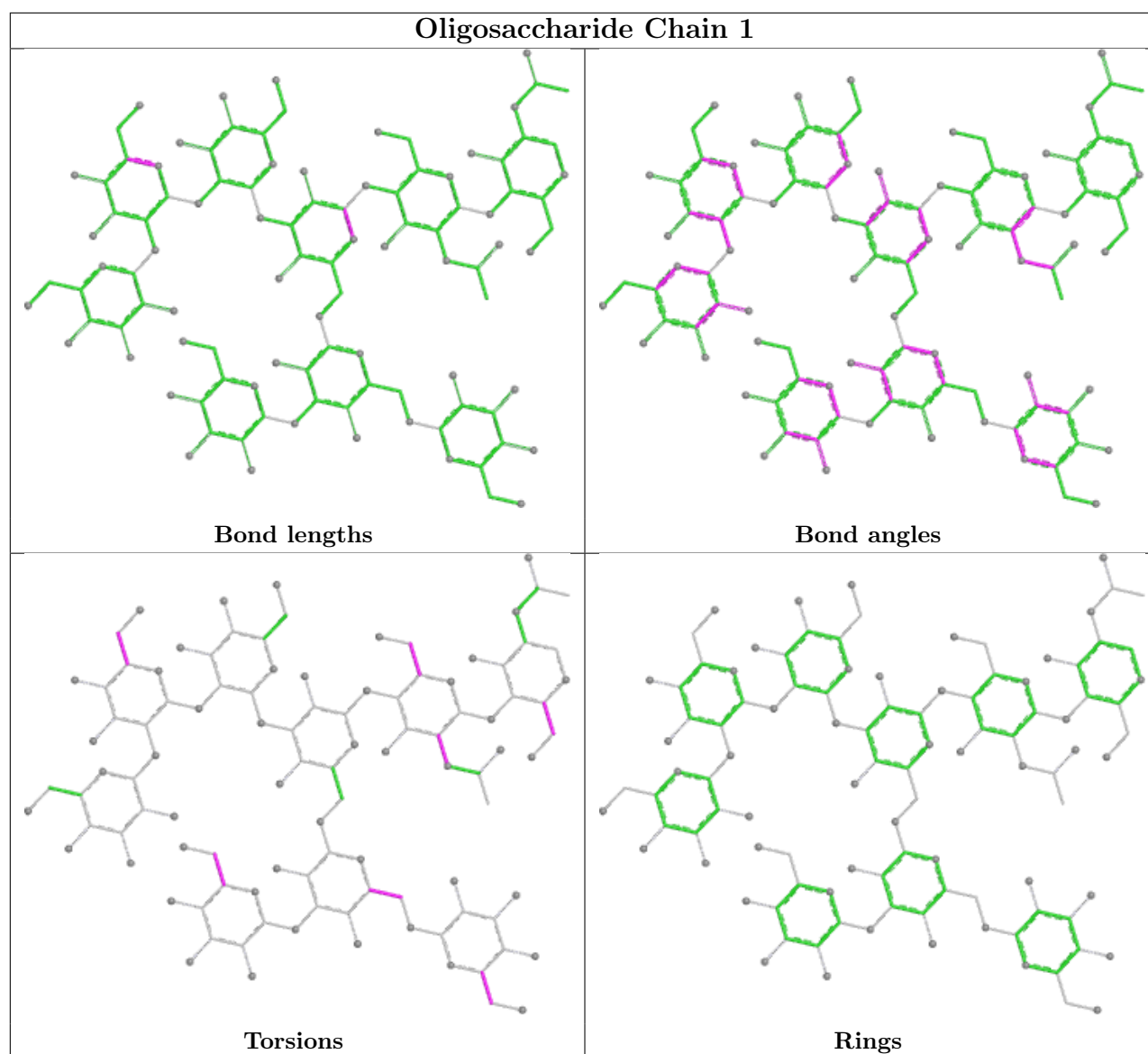












5.6 Ligand geometry [i](#)

9 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	2	611	1	14,14,15	0.27	0	17,19,21	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
14	NAG	c	631	1	14,14,15	0.46	0	17,19,21	0.64	0
14	NAG	C	623	1	14,14,15	0.32	0	17,19,21	0.43	0
14	NAG	c	623	1	14,14,15	0.34	0	17,19,21	0.39	0
14	NAG	c	611	1	14,14,15	0.31	0	17,19,21	0.55	0
14	NAG	2	623	1	14,14,15	0.29	0	17,19,21	0.40	0
14	NAG	2	631	1	14,14,15	0.43	0	17,19,21	0.63	0
14	NAG	C	611	1	14,14,15	0.31	0	17,19,21	0.55	0
14	NAG	C	631	1	14,14,15	0.40	0	17,19,21	0.70	1 (5%)

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	2	611	1	-	0/6/23/26	0/1/1/1
14	NAG	c	631	1	-	2/6/23/26	0/1/1/1
14	NAG	C	623	1	-	2/6/23/26	0/1/1/1
14	NAG	c	623	1	-	2/6/23/26	0/1/1/1
14	NAG	c	611	1	-	0/6/23/26	0/1/1/1
14	NAG	2	623	1	-	2/6/23/26	0/1/1/1
14	NAG	2	631	1	-	2/6/23/26	0/1/1/1
14	NAG	C	611	1	-	0/6/23/26	0/1/1/1
14	NAG	C	631	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	631	NAG	C1-O5-C5	2.26	115.22	112.19

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
14	C	623	NAG	O5-C5-C6-O6
14	c	623	NAG	C4-C5-C6-O6
14	C	623	NAG	C4-C5-C6-O6
14	2	623	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
14	c	623	NAG	O5-C5-C6-O6
14	c	631	NAG	O5-C5-C6-O6
14	2	631	NAG	O5-C5-C6-O6
14	2	623	NAG	O5-C5-C6-O6
14	c	631	NAG	C4-C5-C6-O6
14	2	631	NAG	C4-C5-C6-O6
14	C	631	NAG	O5-C5-C6-O6
14	C	631	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	2	623	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

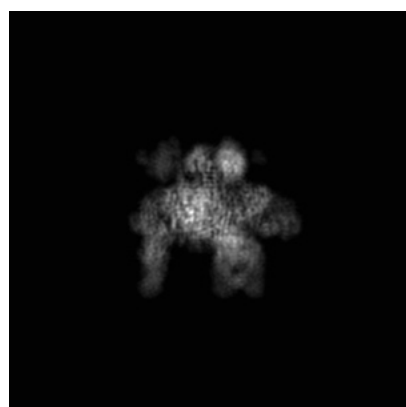
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9320. These allow visual inspection of the internal detail of the map and identification of artifacts.

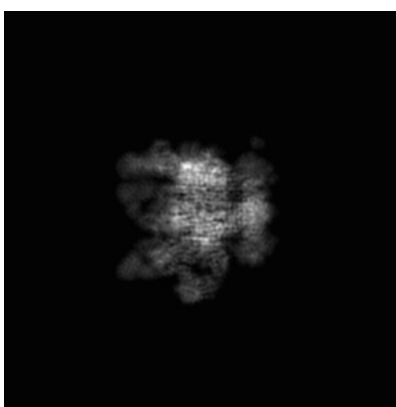
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

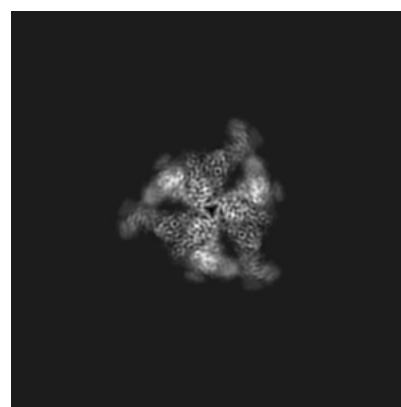
6.1.1 Primary map



X



Y

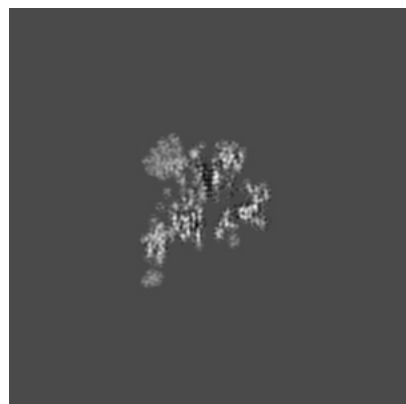


Z

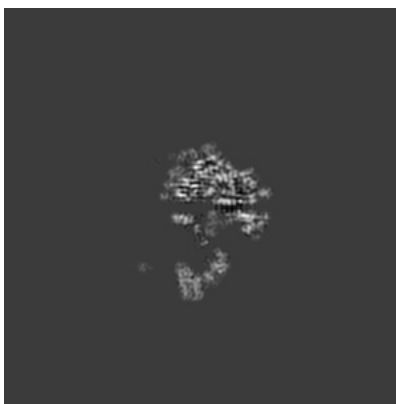
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

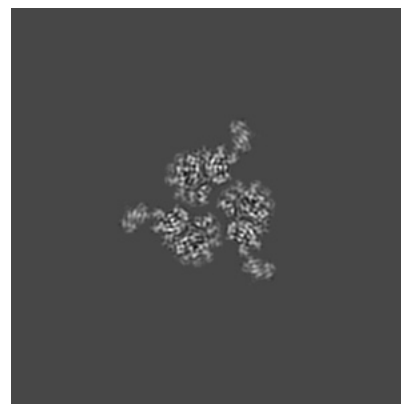
6.2.1 Primary map



X Index: 192



Y Index: 192

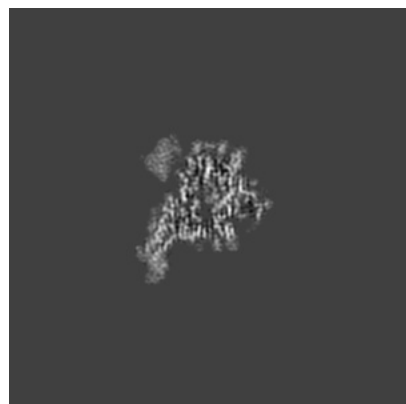


Z Index: 192

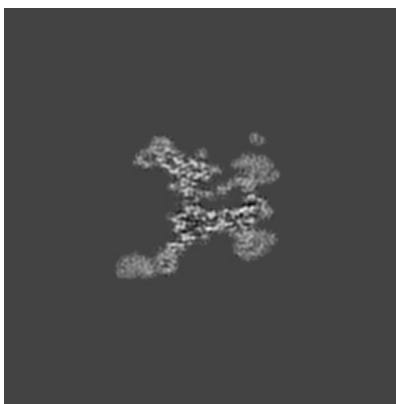
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

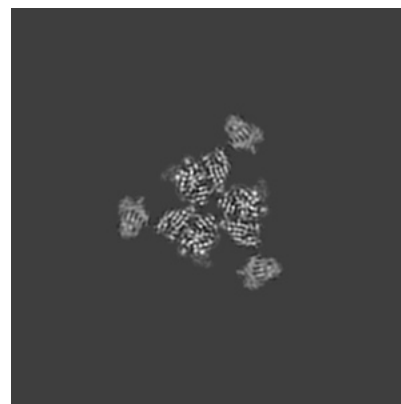
6.3.1 Primary map



X Index: 185



Y Index: 207

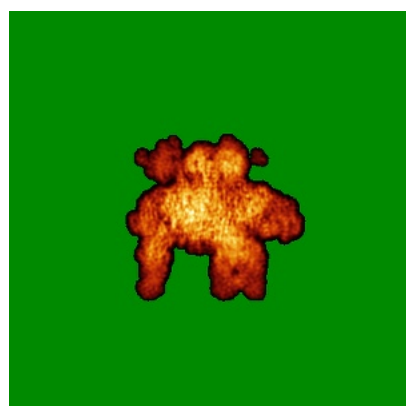


Z Index: 184

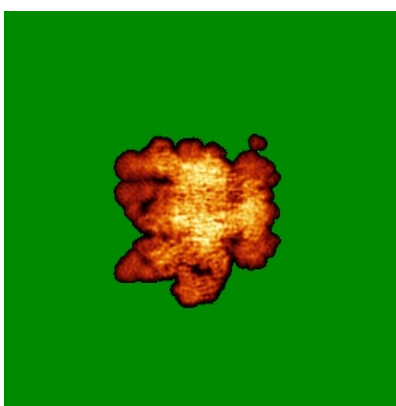
The images above show the largest variance slices of the map in three orthogonal directions.

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

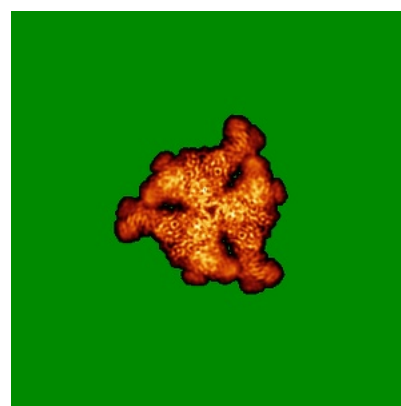
6.4.1 Primary map



X



Y

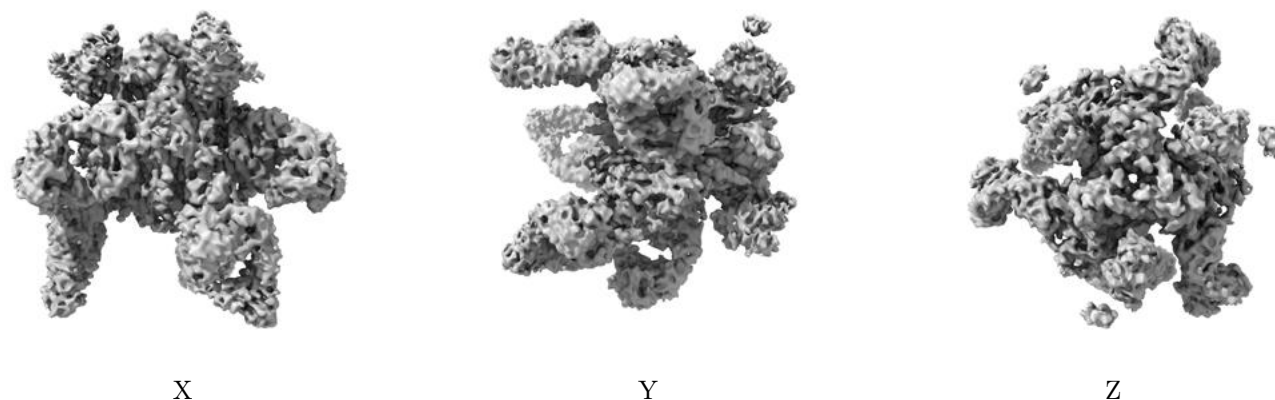


Z

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

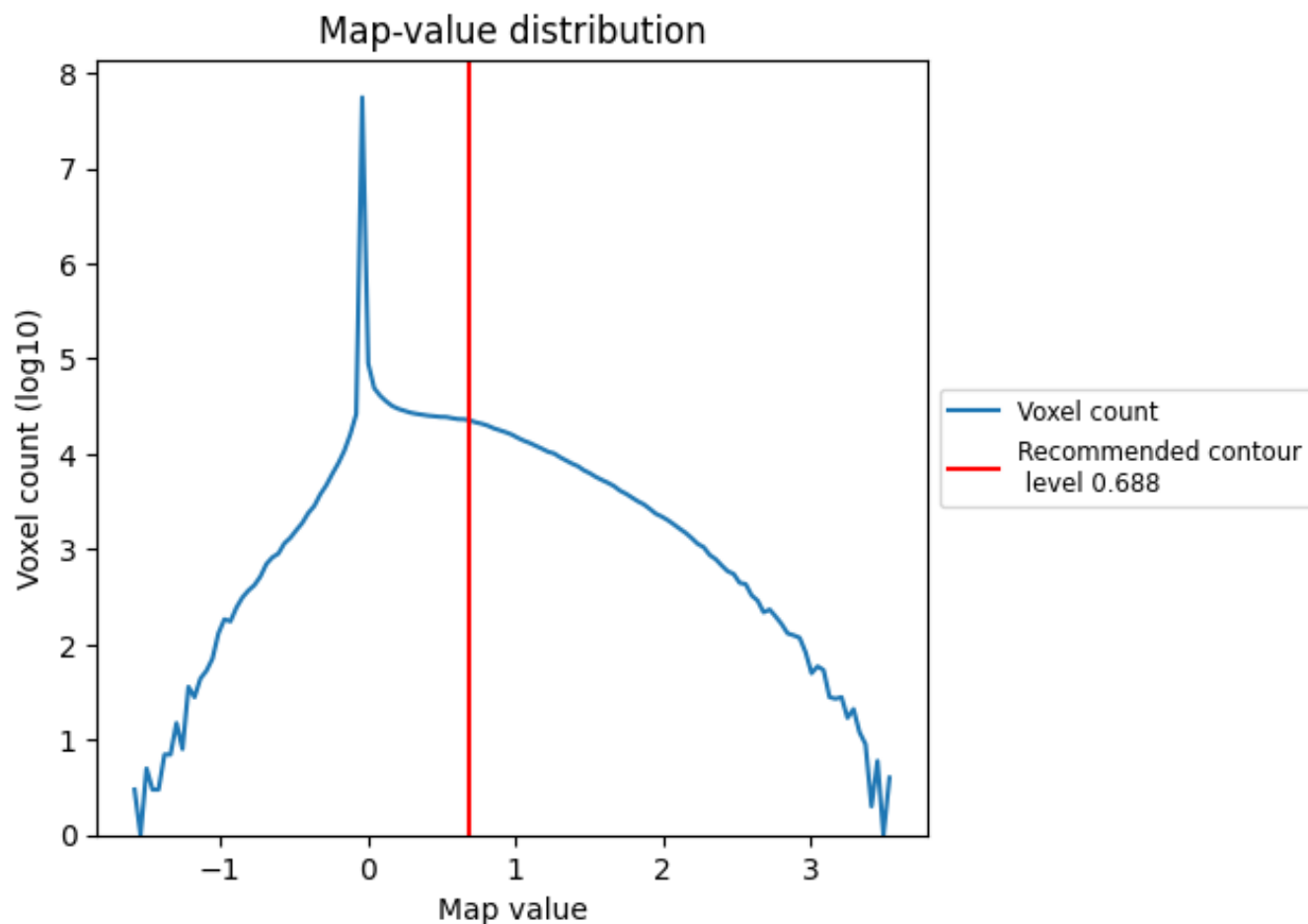
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

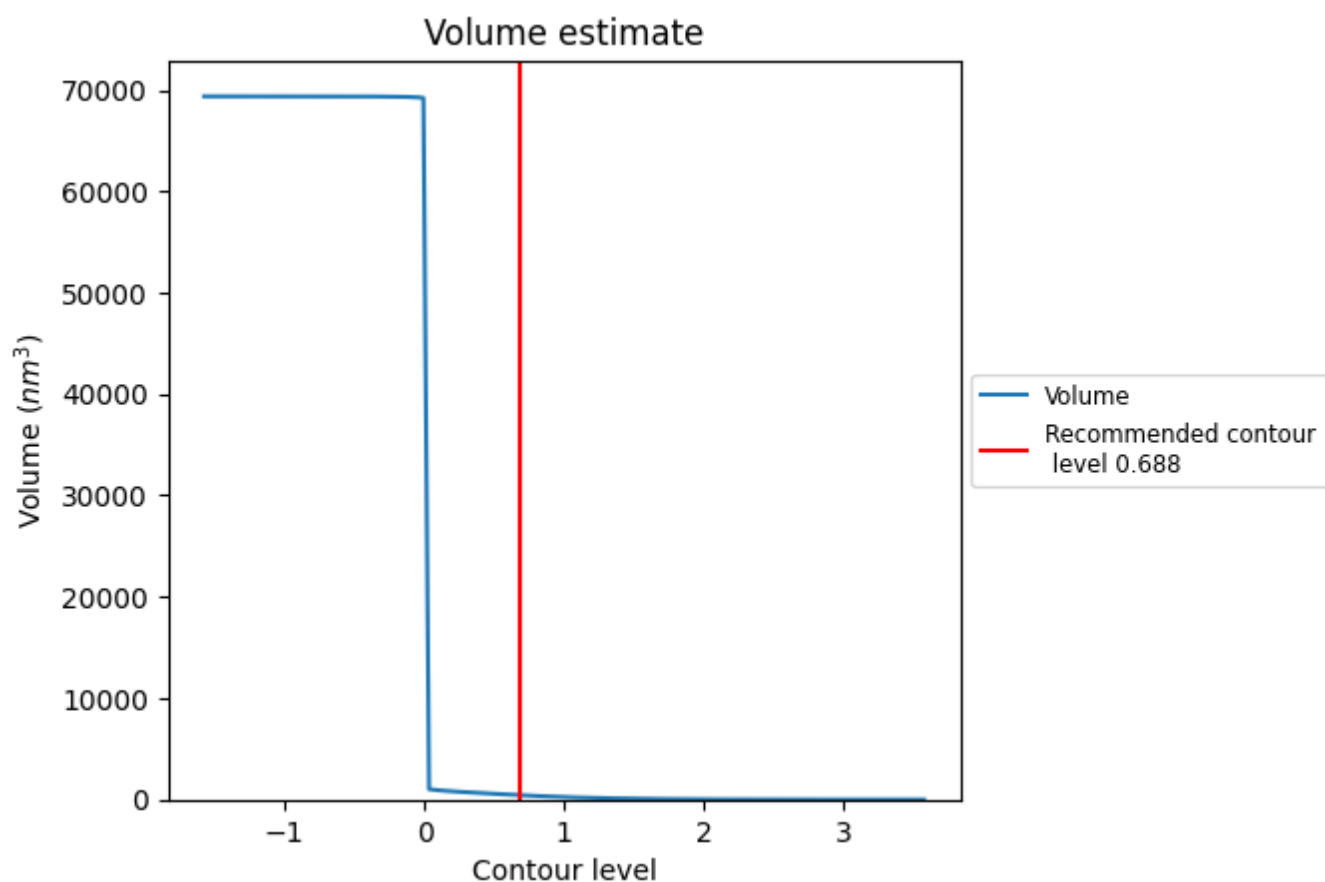
This section contains the results of statistical analysis of the map.

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.

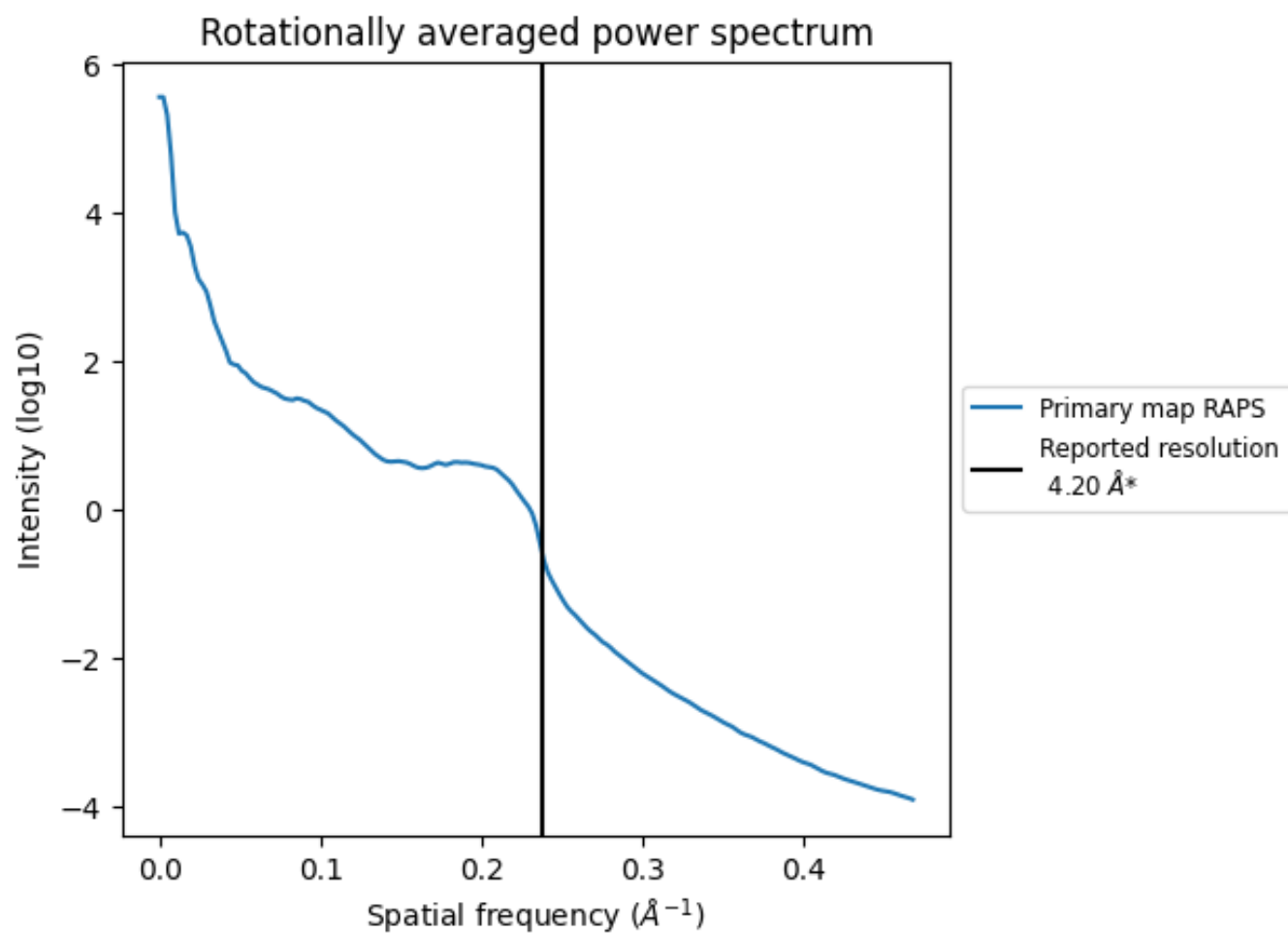
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 431 nm³; this corresponds to an approximate mass of 389 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.238 Å⁻¹

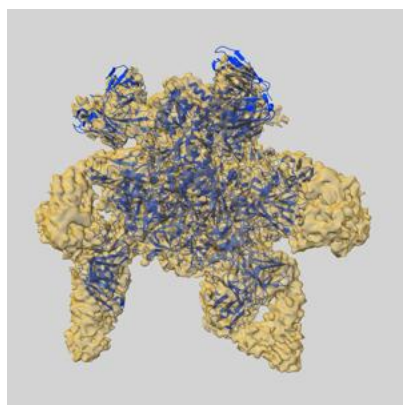
8 Fourier-Shell correlation

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

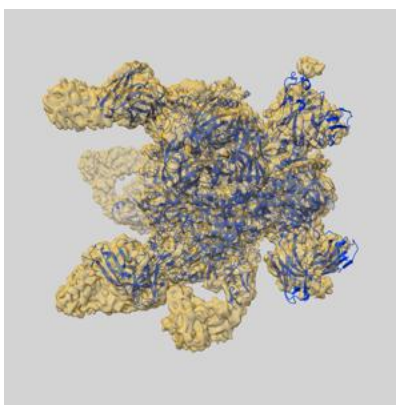
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-9320 and PDB model 6N1W. Per-residue inclusion information can be found in section [3](#) on page [13](#).

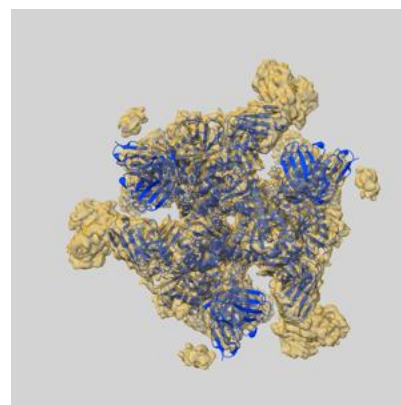
9.1 Map-model overlay [i](#)



X



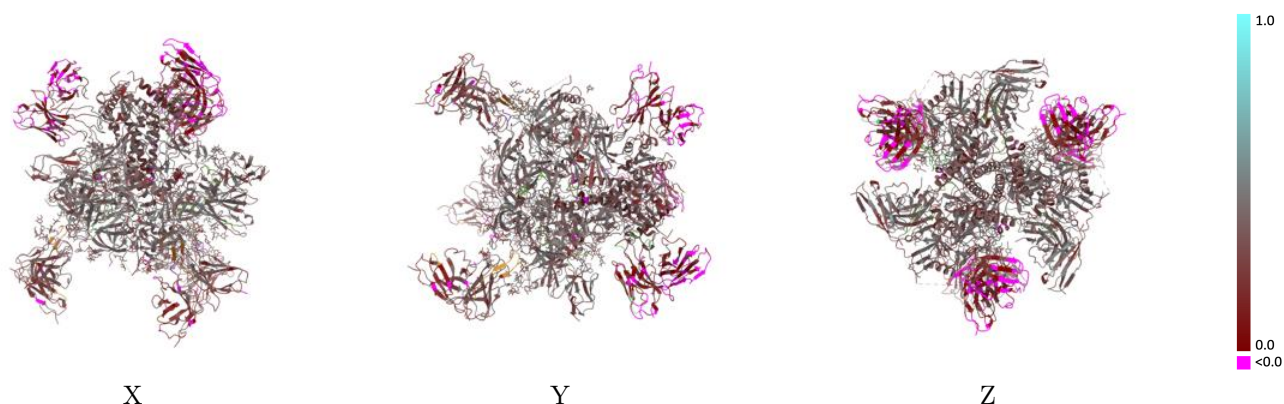
Y



Z

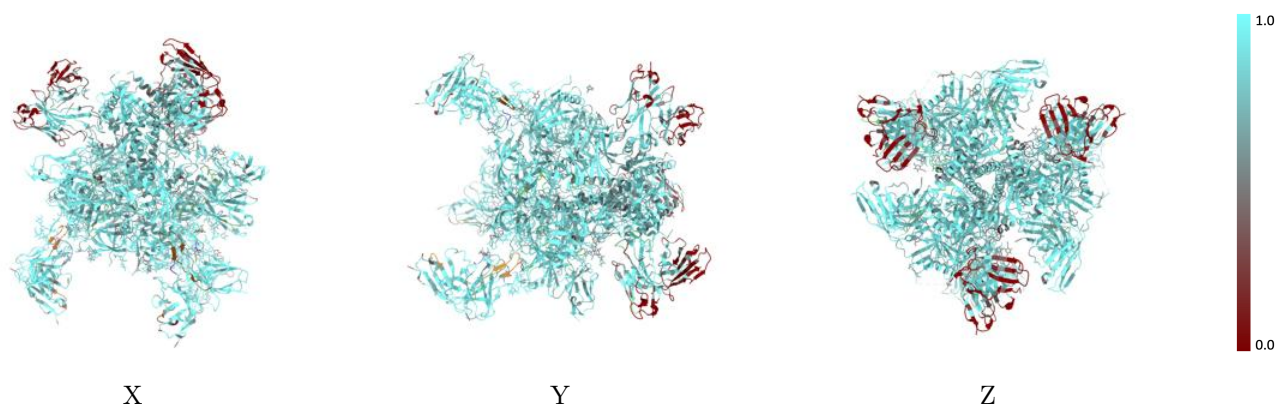
The images above show the 3D surface view of the map at the recommended contour level 0.688 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



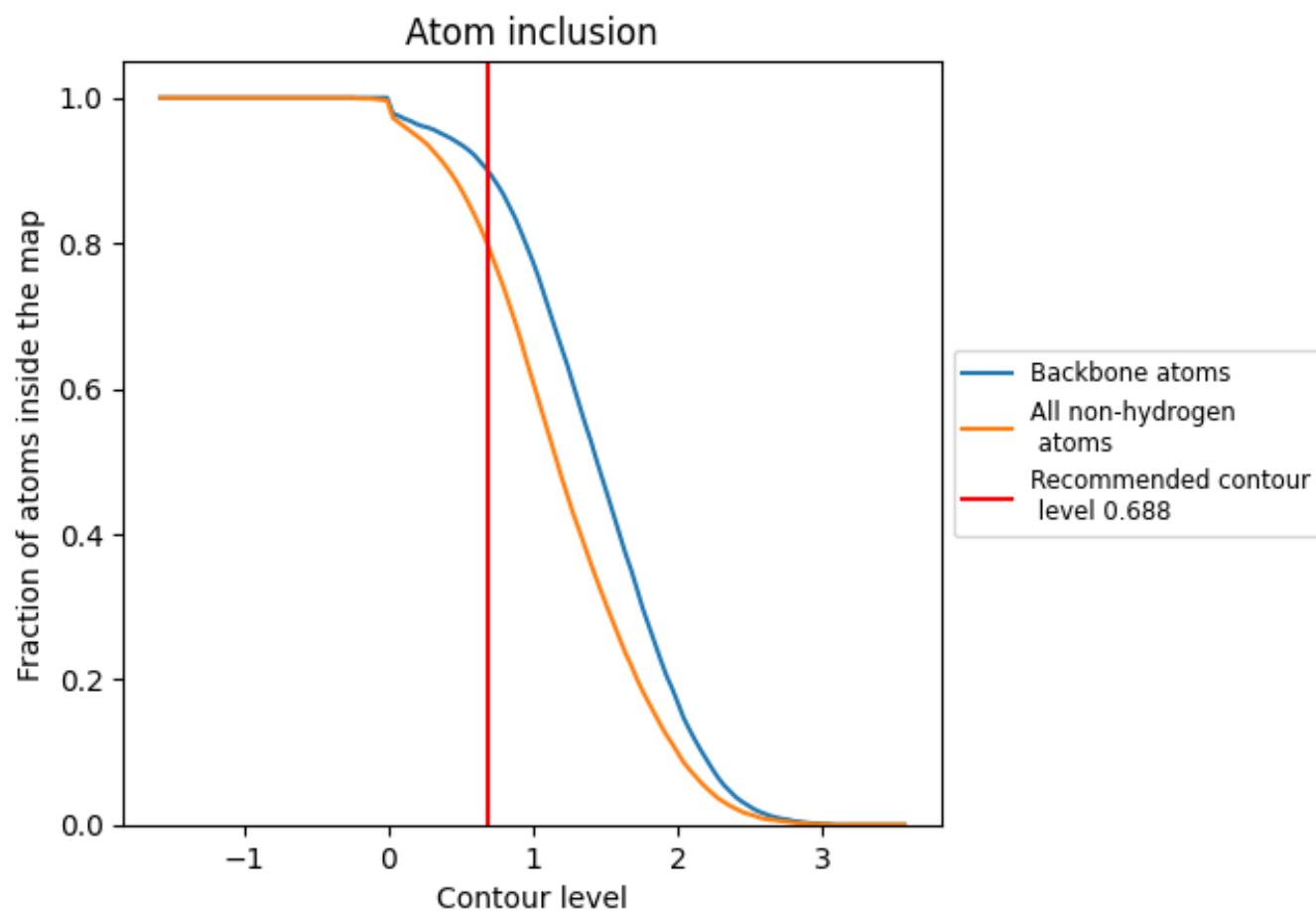
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.688).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 80% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.688) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7980	 0.3140
0	 0.8400	 0.2980
1	 0.9810	 0.3590
2	 0.8520	 0.3760
3	 0.5940	 0.1430
4	 0.4140	 0.0970
5	 0.8490	 0.2770
6	 0.9090	 0.3380
7	 0.8740	 0.3680
8	 0.8620	 0.3990
9	 0.7860	 0.3480
A	 0.8310	 0.3260
AA	 0.7200	 0.3590
B	 0.6430	 0.2560
C	 0.8560	 0.3780
D	 0.8390	 0.3270
E	 0.7400	 0.3140
F	 0.8210	 0.3320
G	 0.8570	 0.3850
H	 0.6020	 0.1440
I	 0.6790	 0.4090
J	 0.7870	 0.3690
K	 0.7500	 0.2990
L	 0.4150	 0.0980
M	 0.8330	 0.2600
N	 0.9080	 0.3280
O	 0.6430	 0.3300
P	 0.8970	 0.4100
Q	 0.8640	 0.3890
R	 0.8770	 0.3600
S	 0.8800	 0.3390
T	 0.9620	 0.3720
U	 0.8570	 0.3570
V	 0.7400	 0.3780
W	 0.6430	 0.2590



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Chain	Atom inclusion	Q-score
X	 0.7800	 0.3180
Y	 0.7860	 0.3330
Z	 0.8930	 0.3870
a	 0.7140	 0.4030
b	 0.7870	 0.3670
c	 0.8500	 0.3710
d	 0.8300	 0.3220
e	 0.7500	 0.3010
f	 0.6430	 0.3270
g	 0.8970	 0.4130
h	 0.5960	 0.1440
i	 0.8600	 0.3690
j	 0.9710	 0.3900
k	 0.7860	 0.3440
l	 0.4130	 0.0930
m	 0.8320	 0.2610
n	 0.9090	 0.3220
o	 0.7400	 0.3670
p	 0.6070	 0.2710
q	 0.8600	 0.3960
r	 0.8760	 0.3700
s	 0.7000	 0.3040
t	 0.7860	 0.3280
u	 0.8570	 0.3780
v	 0.6790	 0.4240
w	 0.7540	 0.3660
x	 0.7860	 0.3400
y	 0.6430	 0.3470
z	 0.8210	 0.3900