



## Full wwPDB EM Validation Report ⓘ

Nov 3, 2025 – 02:22 PM EST

PDB ID : 9OOM / pdb\_00009oom  
EMDB ID : EMD-70666  
Title : Cryo-EM structure of vaccine-elicited antibody T6\_P\_H03 in complex with HIV Env trimer Q23-APEX-GT1  
Authors : Roark, R.S.; Shapiro, L.S.; Kwong, P.D.  
Deposited on : 2025-05-16  
Resolution : 3.50 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev129  
Mogul : 2022.3.0, CSD as543be (2022)  
MolProbity : 4-5-2 with Phenix2.0  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
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) : 2.46

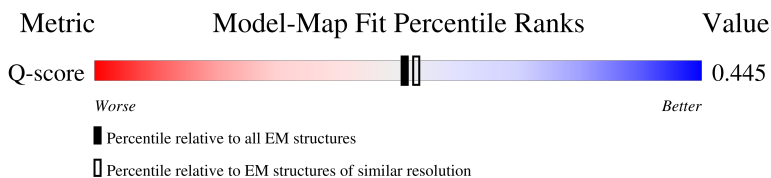
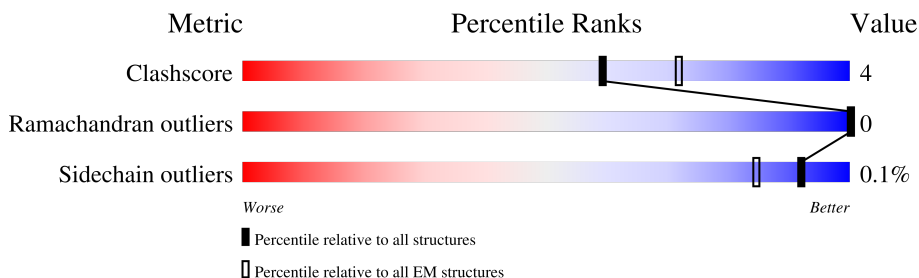
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	210492	15764	-
Ramachandran outliers	207382	16835	-
Sidechain outliers	206894	16415	-
Q-score	-	25397	13950 ( 3.00 - 4.00 )

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	H	235	 5% 50% 6% 44%
1	I	235	 5% 48% 8% 44%
1	J	235	 5% 47% 9% 44%
2	L	214	 5% 38% 12% 50%







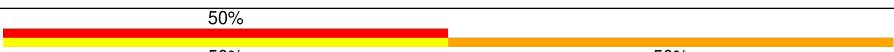



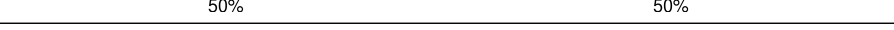
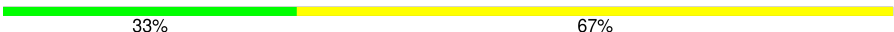




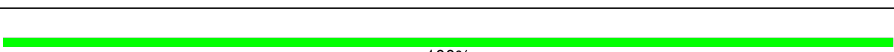
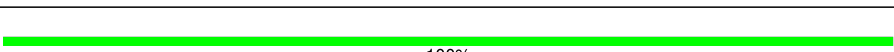

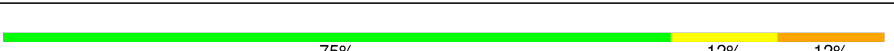


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Mol	Chain	Length	Quality of chain
2	M	214	
2	N	214	
3	b	503	
3	d	503	
3	f	503	
4	c	153	
4	e	153	
4	g	153	
5	A	2	
5	F	2	
5	G	2	
5	K	2	
5	O	2	
5	P	2	
5	Q	2	
5	R	2	
5	S	2	
5	T	2	
5	Y	2	
5	Z	2	
5	a	2	
5	h	2	
5	i	2	
5	j	2	
5	k	2	

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Mol	Chain	Length	Quality of chain
5	l	2	 50% 50%
5	m	2	 50% 50%
5	r	2	 50% 50%
5	s	2	 50% 100%
5	t	2	 50% 50%
5	u	2	 100%
5	v	2	 50% 50%
5	w	2	 50% 50%
5	x	2	 50% 50%
5	y	2	 50% 50%
6	B	3	 33% 67%
6	U	3	 33% 67%
6	n	3	 33% 67%
7	C	5	 100%
7	V	5	 100%
7	o	5	 100%
8	D	5	 80% 60% 40%
9	E	8	 75% 12% 12%
9	q	8	 12% 88% 12%
10	W	5	 80% 60% 40%
10	p	5	 20% 60% 20%
11	X	6	 17% 50% 33%

## 2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 21059 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 T6\_P\_H03 heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	H	132	Total	C	N	O	S	0	0
			1019	643	173	198	5		
1	I	132	Total	C	N	O	S	0	0
			1019	643	173	198	5		
1	J	132	Total	C	N	O	S	0	0
			1019	643	173	198	5		

- Molecule 2 is a protein called T6\_P\_H03 light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	L	107	Total	C	N	O	S	0	0
			810	511	129	167	3		
2	M	107	Total	C	N	O	S	0	0
			810	511	129	167	3		
2	N	107	Total	C	N	O	S	0	0
			810	511	129	167	3		

- Molecule 3 is a protein called HIV Env gp120.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	b	445	Total	C	N	O	S	0	0
			3525	2218	621	657	29		
3	d	445	Total	C	N	O	S	0	0
			3525	2218	621	657	29		
3	f	445	Total	C	N	O	S	0	0
			3525	2218	621	657	29		

- Molecule 4 is a protein called HIV Env gp41 ectodomain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	c	127	Total	C	N	O	S	0	0
			1026	650	184	187	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
4	e	127	Total	C	N	O	S	0	0
			1026	650	184	187	5		
4	g	127	Total	C	N	O	S	0	0
			1026	650	184	187	5		

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



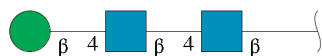
Mol	Chain	Residues	Atoms				AltConf	Trace
5	A	2	Total	C	N	O	0	0
			28	16	2	10		
5	F	2	Total	C	N	O	0	0
			28	16	2	10		
5	G	2	Total	C	N	O	0	0
			28	16	2	10		
5	K	2	Total	C	N	O	0	0
			28	16	2	10		
5	O	2	Total	C	N	O	0	0
			28	16	2	10		
5	P	2	Total	C	N	O	0	0
			28	16	2	10		
5	Q	2	Total	C	N	O	0	0
			28	16	2	10		
5	R	2	Total	C	N	O	0	0
			28	16	2	10		
5	S	2	Total	C	N	O	0	0
			28	16	2	10		
5	T	2	Total	C	N	O	0	0
			28	16	2	10		
5	Y	2	Total	C	N	O	0	0
			28	16	2	10		
5	Z	2	Total	C	N	O	0	0
			28	16	2	10		
5	a	2	Total	C	N	O	0	0
			28	16	2	10		
5	h	2	Total	C	N	O	0	0
			28	16	2	10		
5	i	2	Total	C	N	O	0	0
			28	16	2	10		

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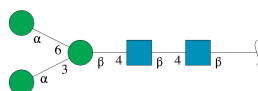
Mol	Chain	Residues	Atoms				AltConf	Trace
5	j	2	Total	C	N	O	0	0
			28	16	2	10		
5	k	2	Total	C	N	O	0	0
			28	16	2	10		
5	l	2	Total	C	N	O	0	0
			28	16	2	10		
5	m	2	Total	C	N	O	0	0
			28	16	2	10		
5	r	2	Total	C	N	O	0	0
			28	16	2	10		
5	s	2	Total	C	N	O	0	0
			28	16	2	10		
5	t	2	Total	C	N	O	0	0
			28	16	2	10		
5	u	2	Total	C	N	O	0	0
			28	16	2	10		
5	v	2	Total	C	N	O	0	0
			28	16	2	10		
5	w	2	Total	C	N	O	0	0
			28	16	2	10		
5	x	2	Total	C	N	O	0	0
			28	16	2	10		
5	y	2	Total	C	N	O	0	0
			28	16	2	10		

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



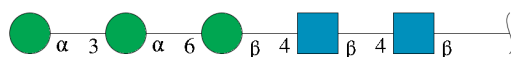
Mol	Chain	Residues	Atoms				AltConf	Trace
6	B	3	Total	C	N	O	0	0
			39	22	2	15		
6	U	3	Total	C	N	O	0	0
			39	22	2	15		
6	n	3	Total	C	N	O	0	0
			39	22	2	15		

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



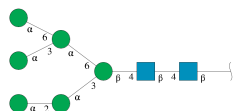
Mol	Chain	Residues	Atoms				AltConf	Trace
7	C	5	Total	C	N	O	0	0
			61	34	2	25		
7	V	5	Total	C	N	O	0	0
			61	34	2	25		
7	o	5	Total	C	N	O	0	0
			61	34	2	25		

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



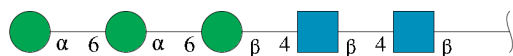
Mol	Chain	Residues	Atoms				AltConf	Trace
8	D	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 9 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



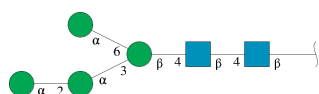
Mol	Chain	Residues	Atoms				AltConf	Trace
9	E	8	Total	C	N	O	0	0
			94	52	2	40		
9	q	8	Total	C	N	O	0	0
			94	52	2	40		

- Molecule 10 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



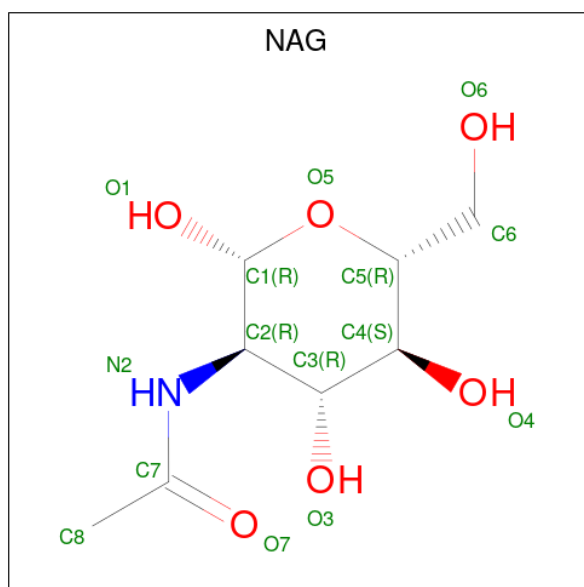
Mol	Chain	Residues	Atoms				AltConf	Trace
10	W	5	Total	C	N	O	0	0
			61	34	2	25		
10	p	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 11 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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

- Molecule 12 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				AltConf
12	b	1	Total	C	N	O	0
			14	8	1	5	
12	b	1	Total	C	N	O	0
			14	8	1	5	
12	b	1	Total	C	N	O	0
			14	8	1	5	
12	b	1	Total	C	N	O	0
			14	8	1	5	
12	b	1	Total	C	N	O	0
			14	8	1	5	
12	b	1	Total	C	N	O	0
			14	8	1	5	
12	c	1	Total	C	N	O	0
			14	8	1	5	
12	c	1	Total	C	N	O	0
			14	8	1	5	
12	c	1	Total	C	N	O	0
			14	8	1	5	
12	d	1	Total	C	N	O	0
			14	8	1	5	
12	d	1	Total	C	N	O	0
			14	8	1	5	
12	d	1	Total	C	N	O	0
			14	8	1	5	
12	d	1	Total	C	N	O	0
			14	8	1	5	
12	d	1	Total	C	N	O	0
			14	8	1	5	
12	d	1	Total	C	N	O	0
			14	8	1	5	
12	e	1	Total	C	N	O	0
			14	8	1	5	
12	e	1	Total	C	N	O	0
			14	8	1	5	
12	e	1	Total	C	N	O	0
			14	8	1	5	
12	f	1	Total	C	N	O	0
			14	8	1	5	
12	f	1	Total	C	N	O	0
			14	8	1	5	

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Mol	Chain	Residues	Atoms				AltConf
12	f	1	Total	C	N	O	0
			14	8	1	5	
12	f	1	Total	C	N	O	0
			14	8	1	5	
12	f	1	Total	C	N	O	0
			14	8	1	5	
12	f	1	Total	C	N	O	0
			14	8	1	5	
12	f	1	Total	C	N	O	0
			14	8	1	5	
12	g	1	Total	C	N	O	0
			14	8	1	5	
12	g	1	Total	C	N	O	0
			14	8	1	5	
12	g	1	Total	C	N	O	0
			14	8	1	5	



[illegible]

Chain M: 

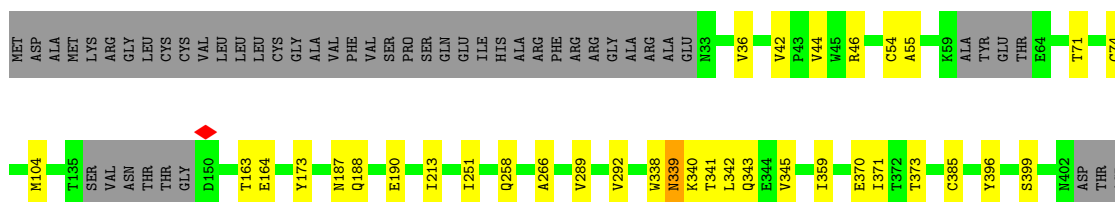
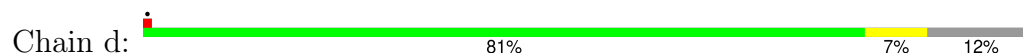
Chain N: 

Chain b:

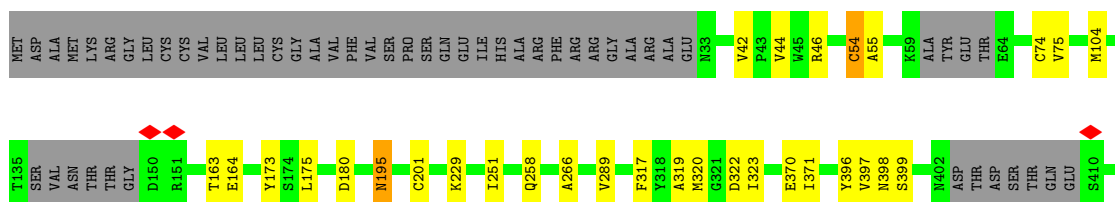
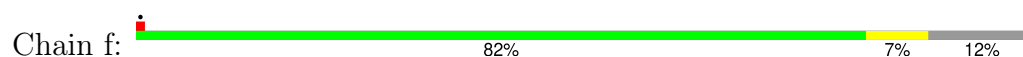
Sequence logo for Chain b. The y-axis represents information content in bits (0.00 to 0.12). The x-axis lists amino acid positions from MET to THR. The logo shows a high conservation of Alanine (ALA) at position 1 (MET) and a mix of residues at position 2 (ASP).



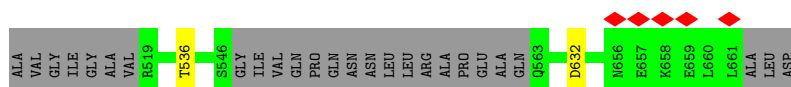
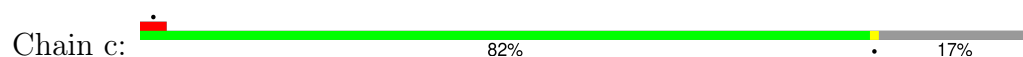
• Molecule 3: HIV Env gp120



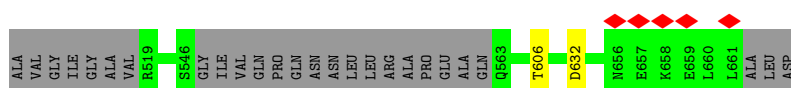
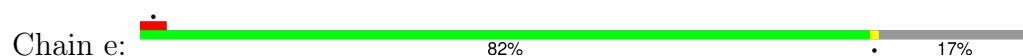
• Molecule 3: HIV Env gp120



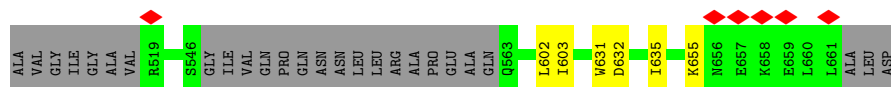
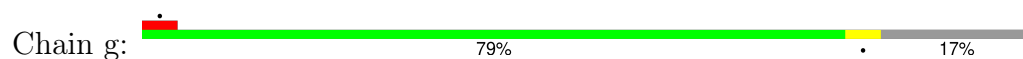
• Molecule 4: HIV Env gp41 ectodomain



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



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



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



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





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

Chain m: A horizontal progress bar for Chain m. It is divided into two equal halves. The left half is green and labeled '50%', and the right half is yellow and labeled '50%'.



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

Chain r: A horizontal progress bar for Chain r. It is divided into two equal halves. The left half is green and labeled '50%', and the right half is yellow and labeled '50%'.



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

Chain s: A horizontal progress bar for Chain s. The first half is red and labeled '50%', and the second half is green and labeled '100%'.



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

Chain t: A horizontal progress bar for Chain t. The first half is red and labeled '50%', and the second half is orange and labeled '50%'.



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

Chain u: A horizontal progress bar for Chain u. It is entirely green and labeled '100%'.



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

Chain v: A horizontal progress bar for Chain v. The first half is red and labeled '50%', and the second half is orange and labeled '50%'.



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



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



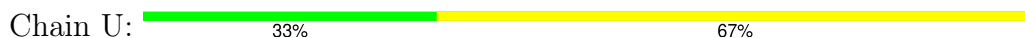
- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



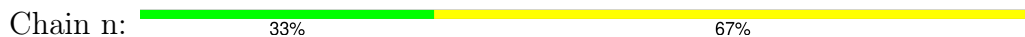
- Molecule 6: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



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





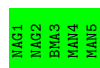
- Molecule 7: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C: 100%



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

Chain V: 100%



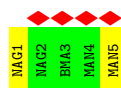
- Molecule 7: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain o: 100%



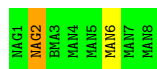
- Molecule 8: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D: 80% 60% 40%



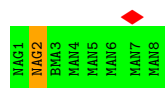
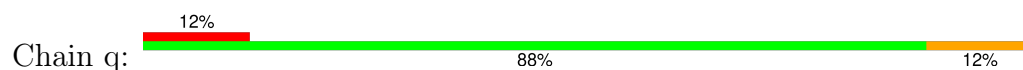
- Molecule 9: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E: 75% 12% 12%

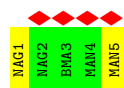
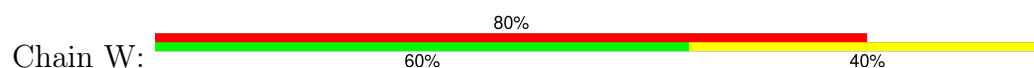


- Molecule 9: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1

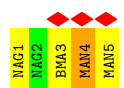
-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



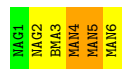
• Molecule 10: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 10: 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



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



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	140657	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	58	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.395	Depositor
Minimum map value	-0.144	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.014	Depositor
Recommended contour level	0.09	Depositor
Map size ( $\text{\AA}$ )	318.72, 318.72, 318.72	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	0.83, 0.83, 0.83	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, MAN, NAG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	H	0.24	0/1042	0.52	0/1408
1	I	0.26	0/1042	0.54	1/1408 (0.1%)
1	J	0.24	0/1042	0.53	1/1408 (0.1%)
2	L	0.61	0/830	0.91	0/1127
2	M	0.61	0/830	0.91	0/1127
2	N	0.61	0/830	0.91	0/1127
3	b	0.28	0/3601	0.50	2/4888 (0.0%)
3	d	0.30	0/3601	0.58	1/4888 (0.0%)
3	f	0.29	0/3601	0.52	2/4888 (0.0%)
4	c	0.19	0/1044	0.43	0/1410
4	e	0.17	0/1044	0.37	0/1410
4	g	0.16	0/1044	0.35	0/1410
All	All	0.33	0/19551	0.58	7/26499 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	d	339	ASN	CB-CA-C	-14.03	88.36	110.81
1	I	90	TYR	CB-CA-C	5.96	117.13	109.80
1	J	90	TYR	CB-CA-C	5.73	116.85	109.80
3	b	195	ASN	CB-CA-C	-5.39	99.86	110.11
3	f	195	ASN	CB-CA-C	-5.38	99.20	110.17
3	f	54	CYS	CB-CA-C	-5.29	100.81	109.53
3	b	54	CYS	CB-CA-C	-5.14	100.00	109.46

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	1019	0	980	12	0
1	I	1019	0	980	14	0
1	J	1019	0	980	16	0
2	L	810	0	778	19	0
2	M	810	0	778	19	0
2	N	810	0	778	20	0
3	b	3525	0	3444	24	0
3	d	3525	0	3444	27	0
3	f	3525	0	3444	23	0
4	c	1026	0	1025	3	0
4	e	1026	0	1025	3	0
4	g	1026	0	1025	5	0
5	A	28	0	25	0	0
5	F	28	0	25	0	0
5	G	28	0	25	1	0
5	K	28	0	25	2	0
5	O	28	0	25	0	0
5	P	28	0	25	1	0
5	Q	28	0	25	0	0
5	R	28	0	25	0	0
5	S	28	0	25	0	0
5	T	28	0	25	0	0
5	Y	28	0	25	0	0
5	Z	28	0	25	0	0
5	a	28	0	25	0	0
5	h	28	0	25	0	0
5	i	28	0	25	0	0
5	j	28	0	25	0	0
5	k	28	0	25	0	0
5	l	28	0	25	0	0
5	m	28	0	25	0	0
5	r	28	0	25	0	0
5	s	28	0	25	0	0
5	t	28	0	25	3	0
5	u	28	0	25	0	0
5	v	28	0	25	1	0
5	w	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	x	28	0	25	0	0
5	y	28	0	25	1	0
6	B	39	0	34	0	0
6	U	39	0	34	0	0
6	n	39	0	34	0	0
7	C	61	0	52	0	0
7	V	61	0	52	0	0
7	o	61	0	52	0	0
8	D	61	0	52	3	0
9	E	94	0	79	2	0
9	q	94	0	79	2	0
10	W	61	0	52	2	0
10	p	61	0	52	6	0
11	X	72	0	61	1	0
12	b	98	0	91	1	0
12	c	42	0	39	0	0
12	d	98	0	91	0	0
12	e	42	0	39	0	0
12	f	98	0	91	1	0
12	g	42	0	39	0	0
All	All	21059	0	20379	184	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (184) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:66:GLY:HA3	2:N:71:PHE:HA	1.81	0.63
2:M:66:GLY:HA3	2:M:71:PHE:HA	1.81	0.62
2:L:61:ARG:HH11	2:L:78:LEU:H	1.47	0.62
5:K:1:NAG:H4	5:K:2:NAG:C7	2.30	0.62
2:L:66:GLY:HA3	2:L:71:PHE:HA	1.81	0.61
5:t:1:NAG:H4	5:t:2:NAG:C7	2.30	0.61
2:N:61:ARG:HH11	2:N:78:LEU:H	1.47	0.60
10:p:3:BMA:H62	10:p:4:MAN:H5	1.84	0.60
2:M:61:ARG:HH11	2:M:78:LEU:H	1.47	0.60
3:d:342:LEU:O	3:d:345:VAL:N	2.35	0.59
1:I:96:ASP:OD1	10:W:5:MAN:H4	2.02	0.59
1:H:55:GLY:HA3	9:E:2:NAG:H82	1.85	0.58
3:d:173:TYR:CE2	10:W:1:NAG:H3	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:f:175:LEU:HD21	10:p:1:NAG:H82	1.85	0.58
2:L:25:ALA:O	2:L:69:THR:HB	2.04	0.58
3:b:336:SER:HA	12:b:603:NAG:H83	1.86	0.58
2:M:25:ALA:O	2:M:69:THR:HB	2.04	0.57
2:N:25:ALA:O	2:N:69:THR:HB	2.04	0.57
3:d:340:LYS:O	3:d:341:THR:C	2.48	0.56
3:b:46:ARG:NH1	4:c:632:ASP:OD2	2.38	0.55
2:L:18:LYS:HG2	2:L:76:SER:HA	1.90	0.54
3:d:258:GLN:NE2	3:d:371:ILE:O	2.41	0.53
3:d:266:ALA:HB3	3:d:289:VAL:HG12	1.90	0.53
3:d:54:CYS:SG	3:d:55:ALA:N	2.82	0.53
2:M:18:LYS:HG2	2:M:76:SER:HA	1.90	0.53
3:b:173:TYR:CE2	8:D:1:NAG:H3	2.44	0.53
4:g:631:TRP:CH2	4:g:635:ILE:HD11	2.44	0.53
3:f:258:GLN:NE2	3:f:371:ILE:O	2.41	0.52
2:N:83:PHE:HD1	2:N:105:GLU:HA	1.74	0.52
2:N:6:GLN:NE2	2:N:86:TYR:O	2.39	0.52
2:N:18:LYS:HG2	2:N:76:SER:HA	1.90	0.52
2:L:30:SER:OG	2:L:31:SER:N	2.43	0.52
3:d:338:TRP:O	3:d:339:ASN:C	2.51	0.52
3:d:359:ILE:O	3:d:467:THR:HA	2.10	0.51
2:M:83:PHE:HD1	2:M:105:GLU:HA	1.74	0.51
3:d:190:GLU:OE1	3:d:190:GLU:N	2.42	0.51
2:L:83:PHE:HD1	2:L:105:GLU:HA	1.74	0.51
3:f:173:TYR:CE2	10:p:1:NAG:H3	2.45	0.51
3:d:373:THR:HB	3:d:385:CYS:O	2.10	0.51
3:b:359:ILE:O	3:b:467:THR:HA	2.10	0.51
2:L:83:PHE:CD1	2:L:105:GLU:HA	2.47	0.50
3:b:46:ARG:NH2	4:c:632:ASP:OD1	2.44	0.50
12:f:607:NAG:H3	5:y:1:NAG:H2	1.93	0.50
3:b:258:GLN:NE2	3:b:371:ILE:O	2.41	0.50
3:f:46:ARG:NH1	4:g:632:ASP:OD2	2.43	0.50
3:f:163:THR:OG1	3:f:164:GLU:N	2.45	0.50
3:d:339:ASN:O	3:d:340:LYS:C	2.55	0.50
3:d:163:THR:OG1	3:d:164:GLU:N	2.45	0.50
2:M:65:SER:O	2:M:72:THR:N	2.39	0.50
2:N:83:PHE:CD1	2:N:105:GLU:HA	2.47	0.50
1:I:6:GLU:N	1:I:6:GLU:OE1	2.45	0.49
2:M:83:PHE:CD1	2:M:105:GLU:HA	2.47	0.49
1:I:33:TRP:HB2	1:I:95:VAL:HG11	1.94	0.49
1:J:6:GLU:OE1	1:J:6:GLU:N	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:f:104:MET:HE1	3:f:251:ILE:HG21	1.95	0.49
4:g:655:LYS:O	4:g:655:LYS:NZ	2.31	0.49
3:f:195:ASN:ND2	3:f:201:CYS:SG	2.85	0.49
2:M:6:GLN:NE2	2:M:86:TYR:O	2.39	0.49
1:H:6:GLU:N	1:H:6:GLU:OE1	2.45	0.49
1:H:29:PHE:O	1:H:71:ARG:NH2	2.45	0.49
1:J:96:ASP:OD2	10:p:5:MAN:H4	2.13	0.49
1:I:29:PHE:O	1:I:71:ARG:NH2	2.45	0.49
1:I:89:VAL:O	1:I:90:TYR:C	2.56	0.49
1:J:33:TRP:HB2	1:J:95:VAL:HG11	1.94	0.49
1:I:89:VAL:HG12	1:I:106:GLY:HA3	1.94	0.49
1:H:33:TRP:HB2	1:H:95:VAL:HG11	1.94	0.49
4:g:602:LEU:C	4:g:603:ILE:HD12	2.38	0.49
1:J:29:PHE:O	1:J:71:ARG:NH2	2.45	0.48
1:H:96:ASP:OD2	8:D:5:MAN:H4	2.12	0.48
3:b:163:THR:OG1	3:b:164:GLU:N	2.45	0.48
3:f:180:ASP:OD2	3:f:422:GLN:N	2.44	0.48
3:b:54:CYS:SG	3:b:55:ALA:N	2.86	0.48
3:f:266:ALA:HB3	3:f:289:VAL:HG12	1.94	0.48
3:d:342:LEU:O	3:d:343:GLN:C	2.55	0.48
3:f:317:PHE:CE2	3:f:319:ALA:HB2	2.49	0.48
1:I:40:THR:HG22	1:I:87:THR:O	2.14	0.48
2:M:22:THR:HB	2:M:72:THR:HG23	1.96	0.47
2:L:22:THR:HB	2:L:72:THR:HG23	1.96	0.47
2:M:18:LYS:HA	2:M:76:SER:HA	1.97	0.47
1:J:13:LYS:HE3	1:J:13:LYS:HA	1.96	0.47
1:H:96:ASP:OD1	8:D:5:MAN:O2	2.31	0.47
2:L:6:GLN:NE2	2:L:86:TYR:O	2.39	0.47
3:b:400:THR:HG22	3:b:402:ASN:H	1.79	0.47
2:N:18:LYS:HA	2:N:76:SER:HA	1.96	0.47
1:H:13:LYS:HE3	1:H:13:LYS:HA	1.96	0.47
1:J:55:GLY:HA3	9:q:2:NAG:H82	1.96	0.47
5:v:1:NAG:H4	5:v:2:NAG:N2	2.30	0.47
1:H:61:ASP:N	1:H:61:ASP:OD1	2.48	0.46
3:d:46:ARG:NH1	4:e:632:ASP:OD2	2.44	0.46
1:J:40:THR:HG22	1:J:87:THR:O	2.16	0.46
2:N:22:THR:HB	2:N:72:THR:HG23	1.96	0.46
2:N:30:SER:OG	2:N:31:SER:N	2.43	0.46
2:L:65:SER:O	2:L:72:THR:N	2.39	0.46
3:f:54:CYS:C	3:f:74:CYS:SG	2.98	0.46
2:L:18:LYS:HA	2:L:76:SER:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:f:322:ASP:OD1	3:f:323:ILE:N	2.48	0.46
1:I:13:LYS:HA	1:I:13:LYS:HE3	1.96	0.46
1:J:96:ASP:CG	10:p:5:MAN:H4	2.41	0.46
3:b:116:LEU:HD11	3:b:434:MET:HE2	1.97	0.46
3:d:71:THR:OG1	3:d:213:ILE:HD11	2.16	0.46
3:d:451:GLY:C	3:d:452:LEU:HD12	2.41	0.46
1:J:89:VAL:O	1:J:90:TYR:C	2.60	0.45
3:b:104:MET:HE1	3:b:251:ILE:HG21	1.99	0.45
3:d:396:TYR:H	3:d:399:SER:HB3	1.81	0.45
1:I:61:ASP:N	1:I:61:ASP:OD1	2.49	0.45
1:J:61:ASP:OD1	1:J:61:ASP:N	2.48	0.45
3:b:54:CYS:C	3:b:74:CYS:SG	3.00	0.45
2:L:61:ARG:HH11	2:L:78:LEU:N	2.13	0.45
3:f:42:VAL:HG13	3:f:44:VAL:HG22	1.98	0.45
3:b:256:SER:OG	3:b:259:LEU:O	2.31	0.44
3:d:42:VAL:HG13	3:d:44:VAL:HG22	1.98	0.44
1:I:11:LEU:HD13	1:I:110:THR:HG23	1.99	0.44
1:I:89:VAL:O	1:I:89:VAL:HG12	2.15	0.44
1:J:11:LEU:HD13	1:J:110:THR:HG23	2.00	0.44
2:N:61:ARG:HH11	2:N:78:LEU:N	2.13	0.44
3:d:187:ASN:ND2	3:d:188:GLN:OE1	2.50	0.44
2:N:68:GLY:O	2:N:71:PHE:CE2	2.71	0.44
5:P:1:NAG:H4	5:P:2:NAG:N2	2.32	0.44
3:b:195:ASN:ND2	3:b:201:CYS:SG	2.91	0.44
3:d:104:MET:HE1	3:d:251:ILE:HG21	1.99	0.44
3:b:401:TRP:O	3:b:402:ASN:C	2.60	0.44
3:f:370:GLU:OE1	3:f:370:GLU:N	2.50	0.44
1:H:11:LEU:HD13	1:H:110:THR:HG23	1.99	0.44
2:M:68:GLY:O	2:M:71:PHE:CE2	2.71	0.43
3:f:320:MET:HE3	3:f:320:MET:HB2	1.77	0.43
3:f:398:ASN:O	3:f:399:SER:C	2.61	0.43
2:L:68:GLY:O	2:L:71:PHE:CE2	2.71	0.43
3:b:370:GLU:OE1	3:b:370:GLU:N	2.50	0.43
3:d:46:ARG:NH2	4:e:632:ASP:OD1	2.51	0.43
1:H:52:ASN:OD1	1:H:54:GLY:N	2.51	0.43
1:H:100(L):LYS:NZ	2:L:92:ASN:OD1	2.41	0.43
1:J:52:ASN:OD1	1:J:54:GLY:N	2.51	0.43
3:b:274:SER:O	5:G:1:NAG:H82	2.18	0.43
1:I:52:ASN:OD1	1:I:54:GLY:N	2.51	0.43
3:b:161:MET:HE2	3:b:309:ILE:HG22	2.01	0.43
3:f:396:TYR:O	3:f:397:VAL:C	2.61	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:f:74:CYS:SG	3:f:75:VAL:N	2.91	0.43
1:J:53:GLY:HA3	9:q:2:NAG:H2	2.01	0.43
2:N:65:SER:O	2:N:72:THR:N	2.39	0.43
1:J:100(L):LYS:NZ	2:N:92:ASN:OD1	2.41	0.42
3:f:46:ARG:NH2	4:g:632:ASP:OD1	2.52	0.42
2:N:75:ILE:HG22	2:N:77:SER:H	1.84	0.42
2:M:61:ARG:HH11	2:M:78:LEU:N	2.13	0.42
2:M:80:PRO:HA	2:M:83:PHE:HE2	1.84	0.42
2:N:80:PRO:HA	2:N:83:PHE:HE2	1.84	0.42
2:L:61:ARG:CZ	2:L:79:GLN:HG2	2.49	0.42
2:M:61:ARG:CZ	2:M:79:GLN:HG2	2.49	0.42
2:M:75:ILE:HG22	2:M:77:SER:H	1.84	0.42
2:L:4:MET:HE1	2:L:25:ALA:HB1	2.02	0.42
3:d:54:CYS:C	3:d:74:CYS:SG	3.02	0.42
3:d:338:TRP:O	3:d:341:THR:OG1	2.37	0.42
3:d:370:GLU:OE1	3:d:370:GLU:N	2.50	0.42
1:J:68:ILE:O	1:J:81:GLN:N	2.52	0.42
3:b:255:VAL:HG11	3:b:426:MET:HE2	2.01	0.42
2:M:85:THR:HA	2:M:102:THR:O	2.20	0.42
3:f:323:ILE:HD12	3:f:323:ILE:HA	1.87	0.42
2:N:61:ARG:CZ	2:N:79:GLN:HG2	2.49	0.42
3:b:42:VAL:HG13	3:b:44:VAL:HG22	2.01	0.41
1:I:100(L):LYS:NZ	2:M:92:ASN:OD1	2.41	0.41
1:H:53:GLY:HA3	9:E:2:NAG:H2	2.02	0.41
3:b:398:ASN:O	3:b:399:SER:C	2.63	0.41
2:M:4:MET:HE1	2:M:25:ALA:HB1	2.02	0.41
2:N:4:MET:HE1	2:N:25:ALA:HB1	2.02	0.41
2:N:85:THR:HA	2:N:102:THR:O	2.20	0.41
2:L:75:ILE:HG22	2:L:77:SER:H	1.84	0.41
2:M:30:SER:OG	2:M:31:SER:N	2.43	0.41
5:t:1:NAG:H4	5:t:2:NAG:N2	2.35	0.41
3:d:36:VAL:O	4:e:606:THR:OG1	2.39	0.41
2:N:67:SER:O	2:N:68:GLY:C	2.64	0.41
2:L:80:PRO:HA	2:L:83:PHE:HE2	1.84	0.41
1:J:96:ASP:OD1	10:p:5:MAN:H4	2.20	0.41
3:b:184:ILE:HD12	3:b:190:GLU:O	2.20	0.41
1:I:38:ARG:HB2	1:I:88:ALA:CB	2.51	0.41
3:b:67:ASN:OD1	3:b:70:ALA:N	2.54	0.41
5:K:1:NAG:H4	5:K:2:NAG:N2	2.35	0.41
2:L:85:THR:HA	2:L:102:THR:O	2.20	0.40
3:f:54:CYS:SG	3:f:55:ALA:N	2.93	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:X:4:MAN:H2	11:X:5:MAN:H2	1.81	0.40
4:c:536:THR:O	4:c:536:THR:OG1	2.36	0.40
3:d:292:VAL:HG21	3:d:341:THR:HG21	2.04	0.40
3:d:340:LYS:HE3	3:d:340:LYS:HB2	1.81	0.40
3:f:229:LYS:HD3	5:t:1:NAG:H61	2.04	0.40
3:b:350:ARG:NH2	3:b:396:TYR:O	2.55	0.40
3:f:457:ASP:OD2	3:f:467:THR:OG1	2.23	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	130/235 (55%)	120 (92%)	10 (8%)	0	100	100
1	I	130/235 (55%)	118 (91%)	12 (9%)	0	100	100
1	J	130/235 (55%)	118 (91%)	12 (9%)	0	100	100
2	L	105/214 (49%)	98 (93%)	7 (7%)	0	100	100
2	M	105/214 (49%)	98 (93%)	7 (7%)	0	100	100
2	N	105/214 (49%)	98 (93%)	7 (7%)	0	100	100
3	b	437/503 (87%)	420 (96%)	17 (4%)	0	100	100
3	d	437/503 (87%)	414 (95%)	23 (5%)	0	100	100
3	f	437/503 (87%)	417 (95%)	20 (5%)	0	100	100
4	c	123/153 (80%)	120 (98%)	3 (2%)	0	100	100
4	e	123/153 (80%)	118 (96%)	5 (4%)	0	100	100
4	g	123/153 (80%)	120 (98%)	3 (2%)	0	100	100
All	All	2385/3315 (72%)	2259 (95%)	126 (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	H	108/198 (54%)	108 (100%)	0	100	100
1	I	108/198 (54%)	108 (100%)	0	100	100
1	J	108/198 (54%)	108 (100%)	0	100	100
2	L	92/191 (48%)	92 (100%)	0	100	100
2	M	92/191 (48%)	92 (100%)	0	100	100
2	N	92/191 (48%)	92 (100%)	0	100	100
3	b	398/446 (89%)	397 (100%)	1 (0%)	91	96
3	d	398/446 (89%)	398 (100%)	0	100	100
3	f	398/446 (89%)	397 (100%)	1 (0%)	91	96
4	c	110/128 (86%)	110 (100%)	0	100	100
4	e	110/128 (86%)	110 (100%)	0	100	100
4	g	110/128 (86%)	110 (100%)	0	100	100
All	All	2124/2889 (74%)	2122 (100%)	2 (0%)	92	97

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	b	445	CYS
3	f	445	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
4	c	590	GLN
3	d	185	ASN
3	d	195	ASN
3	f	188	GLN
4	g	590	GLN
4	g	652	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

115 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	$\# Z  > 2$
5	NAG	A	1	3,5	14,14,15	0.69	0	17,19,21	0.88	1 (5%)
5	NAG	A	2	5	14,14,15	0.68	0	17,19,21	0.88	0
6	NAG	B	1	3,6	14,14,15	0.69	0	17,19,21	0.82	0
6	NAG	B	2	6	14,14,15	0.69	0	17,19,21	0.93	1 (5%)
6	BMA	B	3	6	11,11,12	0.82	0	15,15,17	2.17	4 (26%)
7	NAG	C	1	3,7	14,14,15	0.40	0	17,19,21	0.59	0
7	NAG	C	2	7	14,14,15	0.39	0	17,19,21	0.59	0
7	BMA	C	3	7	11,11,12	0.40	0	15,15,17	0.46	0
7	MAN	C	4	7	11,11,12	0.34	0	15,15,17	0.50	0
7	MAN	C	5	7	11,11,12	0.29	0	15,15,17	0.52	0
8	NAG	D	1	3,8	14,14,15	0.38	0	17,19,21	0.68	0
8	NAG	D	2	8	14,14,15	0.38	0	17,19,21	0.62	0
8	BMA	D	3	8	11,11,12	0.33	0	15,15,17	0.51	0
8	MAN	D	4	8	11,11,12	0.34	0	15,15,17	0.52	0
8	MAN	D	5	8	11,11,12	0.44	0	15,15,17	0.58	0
9	NAG	E	1	3,9	14,14,15	0.42	0	17,19,21	0.66	0
9	NAG	E	2	9	14,14,15	0.38	0	17,19,21	0.80	1 (5%)
9	BMA	E	3	9	11,11,12	0.58	0	15,15,17	0.51	0
9	MAN	E	4	9	11,11,12	0.41	0	15,15,17	0.51	0
9	MAN	E	5	9	11,11,12	0.32	0	15,15,17	0.57	0
9	MAN	E	6	9	11,11,12	0.43	0	15,15,17	0.88	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
9	MAN	E	7	9	11,11,12	0.32	0	15,15,17	0.51	0
9	MAN	E	8	9	11,11,12	0.31	0	15,15,17	0.54	0
5	NAG	F	1	3,5	14,14,15	0.36	0	17,19,21	0.76	0
5	NAG	F	2	5	14,14,15	0.36	0	17,19,21	0.67	0
5	NAG	G	1	3,5	14,14,15	0.38	0	17,19,21	0.73	0
5	NAG	G	2	5	14,14,15	0.42	0	17,19,21	0.77	1 (5%)
5	NAG	K	1	3,5	14,14,15	0.38	0	17,19,21	0.58	0
5	NAG	K	2	5	14,14,15	0.40	0	17,19,21	0.91	2 (11%)
5	NAG	O	1	3,5	14,14,15	0.38	0	17,19,21	0.81	0
5	NAG	O	2	5	14,14,15	0.39	0	17,19,21	0.54	0
5	NAG	P	1	3,5	14,14,15	0.39	0	17,19,21	0.63	0
5	NAG	P	2	5	14,14,15	0.40	0	17,19,21	0.90	1 (5%)
5	NAG	Q	1	3,5	14,14,15	0.37	0	17,19,21	0.67	0
5	NAG	Q	2	5	14,14,15	0.39	0	17,19,21	0.82	1 (5%)
5	NAG	R	1	3,5	14,14,15	0.38	0	17,19,21	0.63	0
5	NAG	R	2	5	14,14,15	0.37	0	17,19,21	0.64	1 (5%)
5	NAG	S	1	3,5	14,14,15	0.38	0	17,19,21	0.59	0
5	NAG	S	2	5	14,14,15	0.37	0	17,19,21	0.83	1 (5%)
5	NAG	T	1	3,5	14,14,15	0.68	0	17,19,21	0.87	1 (5%)
5	NAG	T	2	5	14,14,15	0.68	0	17,19,21	0.88	0
6	NAG	U	1	3,6	14,14,15	0.69	0	17,19,21	0.81	0
6	NAG	U	2	6	14,14,15	0.70	0	17,19,21	0.93	1 (5%)
6	BMA	U	3	6	11,11,12	0.80	0	15,15,17	2.19	4 (26%)
7	NAG	V	1	3,7	14,14,15	0.41	0	17,19,21	0.56	0
7	NAG	V	2	7	14,14,15	0.36	0	17,19,21	0.53	0
7	BMA	V	3	7	11,11,12	0.31	0	15,15,17	0.50	0
7	MAN	V	4	7	11,11,12	0.32	0	15,15,17	0.54	0
7	MAN	V	5	7	11,11,12	0.30	0	15,15,17	0.52	0
10	NAG	W	1	3,10	14,14,15	0.38	0	17,19,21	0.72	0
10	NAG	W	2	10	14,14,15	0.39	0	17,19,21	0.71	0
10	BMA	W	3	10	11,11,12	0.31	0	15,15,17	0.45	0
10	MAN	W	4	10	11,11,12	0.29	0	15,15,17	0.51	0
10	MAN	W	5	10	11,11,12	0.33	0	15,15,17	0.60	0
11	NAG	X	1	11,3	14,14,15	0.72	0	17,19,21	0.92	0
11	NAG	X	2	11	14,14,15	0.67	0	17,19,21	1.11	1 (5%)
11	BMA	X	3	11	11,11,12	0.84	0	15,15,17	2.35	5 (33%)
11	MAN	X	4	11	11,11,12	0.68	0	15,15,17	1.11	1 (6%)
11	MAN	X	5	11	11,11,12	0.74	0	15,15,17	0.97	1 (6%)
11	MAN	X	6	11	11,11,12	0.74	0	15,15,17	0.93	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	NAG	Y	1	3,5	14,14,15	0.69	0	17,19,21	1.04	1 (5%)
5	NAG	Y	2	5	14,14,15	0.68	0	17,19,21	0.99	1 (5%)
5	NAG	Z	1	3,5	14,14,15	0.68	0	17,19,21	0.97	1 (5%)
5	NAG	Z	2	5	14,14,15	0.68	0	17,19,21	0.84	0
5	NAG	a	1	3,5	14,14,15	0.72	0	17,19,21	0.96	0
5	NAG	a	2	5	14,14,15	0.80	0	17,19,21	1.32	2 (11%)
5	NAG	h	1	3,5	14,14,15	0.69	0	17,19,21	1.11	2 (11%)
5	NAG	h	2	5	14,14,15	0.76	0	17,19,21	0.91	1 (5%)
5	NAG	i	1	3,5	14,14,15	0.71	0	17,19,21	0.94	1 (5%)
5	NAG	i	2	5	14,14,15	0.73	0	17,19,21	1.29	2 (11%)
5	NAG	j	1	3,5	14,14,15	0.72	0	17,19,21	1.01	1 (5%)
5	NAG	j	2	5	14,14,15	0.76	0	17,19,21	1.23	1 (5%)
5	NAG	k	1	3,5	14,14,15	0.40	0	17,19,21	0.67	0
5	NAG	k	2	5	14,14,15	0.39	0	17,19,21	0.70	1 (5%)
5	NAG	l	1	3,5	14,14,15	0.39	0	17,19,21	0.54	0
5	NAG	l	2	5	14,14,15	0.40	0	17,19,21	1.39	2 (11%)
5	NAG	m	1	3,5	14,14,15	0.70	0	17,19,21	0.92	1 (5%)
5	NAG	m	2	5	14,14,15	0.68	0	17,19,21	0.88	0
6	NAG	n	1	3,6	14,14,15	0.69	0	17,19,21	0.83	0
6	NAG	n	2	6	14,14,15	0.71	0	17,19,21	0.95	2 (11%)
6	BMA	n	3	6	11,11,12	0.82	0	15,15,17	2.19	4 (26%)
7	NAG	o	1	3,7	14,14,15	0.41	0	17,19,21	0.56	0
7	NAG	o	2	7	14,14,15	0.40	0	17,19,21	0.62	0
7	BMA	o	3	7	11,11,12	0.37	0	15,15,17	0.49	0
7	MAN	o	4	7	11,11,12	0.32	0	15,15,17	0.57	0
7	MAN	o	5	7	11,11,12	0.30	0	15,15,17	0.52	0
10	NAG	p	1	3,10	14,14,15	0.39	0	17,19,21	0.76	0
10	NAG	p	2	10	14,14,15	0.37	0	17,19,21	0.75	0
10	BMA	p	3	10	11,11,12	0.34	0	15,15,17	0.55	0
10	MAN	p	4	10	11,11,12	0.28	0	15,15,17	0.89	1 (6%)
10	MAN	p	5	10	11,11,12	0.31	0	15,15,17	0.57	0
9	NAG	q	1	3,9	14,14,15	0.43	0	17,19,21	0.55	0
9	NAG	q	2	9	14,14,15	0.36	0	17,19,21	1.16	1 (5%)
9	BMA	q	3	9	11,11,12	0.52	0	15,15,17	0.57	0
9	MAN	q	4	9	11,11,12	0.39	0	15,15,17	0.61	0
9	MAN	q	5	9	11,11,12	0.38	0	15,15,17	0.54	0
9	MAN	q	6	9	11,11,12	0.46	0	15,15,17	0.54	0
9	MAN	q	7	9	11,11,12	0.30	0	15,15,17	0.49	0
9	MAN	q	8	9	11,11,12	0.37	0	15,15,17	0.52	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	NAG	r	1	3,5	14,14,15	0.37	0	17,19,21	0.75	0
5	NAG	r	2	5	14,14,15	0.38	0	17,19,21	0.67	1 (5%)
5	NAG	s	1	3,5	14,14,15	0.40	0	17,19,21	0.72	0
5	NAG	s	2	5	14,14,15	0.38	0	17,19,21	0.44	0
5	NAG	t	1	3,5	14,14,15	0.38	0	17,19,21	0.56	0
5	NAG	t	2	5	14,14,15	0.40	0	17,19,21	0.91	2 (11%)
5	NAG	u	1	3,5	14,14,15	0.40	0	17,19,21	0.76	0
5	NAG	u	2	5	14,14,15	0.38	0	17,19,21	0.55	0
5	NAG	v	1	3,5	14,14,15	0.39	0	17,19,21	0.60	0
5	NAG	v	2	5	14,14,15	0.39	0	17,19,21	0.91	1 (5%)
5	NAG	w	1	3,5	14,14,15	0.39	0	17,19,21	0.68	0
5	NAG	w	2	5	14,14,15	0.38	0	17,19,21	0.82	1 (5%)
5	NAG	x	1	3,5	14,14,15	0.39	0	17,19,21	0.67	0
5	NAG	x	2	5	14,14,15	0.37	0	17,19,21	0.64	1 (5%)
5	NAG	y	1	3,5	14,14,15	0.36	0	17,19,21	0.71	1 (5%)
5	NAG	y	2	5	14,14,15	0.37	0	17,19,21	0.82	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
5	NAG	A	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	A	2	5	-	0/6/23/26	0/1/1/1
6	NAG	B	1	3,6	-	0/6/23/26	0/1/1/1
6	NAG	B	2	6	-	2/6/23/26	0/1/1/1
6	BMA	B	3	6	-	0/2/19/22	0/1/1/1
7	NAG	C	1	3,7	-	0/6/23/26	0/1/1/1
7	NAG	C	2	7	-	0/6/23/26	0/1/1/1
7	BMA	C	3	7	-	0/2/19/22	0/1/1/1
7	MAN	C	4	7	-	0/2/19/22	0/1/1/1
7	MAN	C	5	7	-	0/2/19/22	0/1/1/1
8	NAG	D	1	3,8	-	0/6/23/26	0/1/1/1
8	NAG	D	2	8	-	0/6/23/26	0/1/1/1
8	BMA	D	3	8	-	0/2/19/22	0/1/1/1
8	MAN	D	4	8	-	0/2/19/22	0/1/1/1
8	MAN	D	5	8	-	1/2/19/22	0/1/1/1
9	NAG	E	1	3,9	-	0/6/23/26	0/1/1/1
9	NAG	E	2	9	-	2/6/23/26	0/1/1/1
9	BMA	E	3	9	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MAN	E	4	9	-	0/2/19/22	0/1/1/1
9	MAN	E	5	9	-	0/2/19/22	0/1/1/1
9	MAN	E	6	9	-	0/2/19/22	0/1/1/1
9	MAN	E	7	9	-	0/2/19/22	0/1/1/1
9	MAN	E	8	9	-	0/2/19/22	0/1/1/1
5	NAG	F	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	F	2	5	-	2/6/23/26	0/1/1/1
5	NAG	G	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	G	2	5	-	1/6/23/26	0/1/1/1
5	NAG	K	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	K	2	5	-	0/6/23/26	0/1/1/1
5	NAG	O	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	O	2	5	-	1/6/23/26	0/1/1/1
5	NAG	P	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	P	2	5	-	2/6/23/26	0/1/1/1
5	NAG	Q	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	Q	2	5	-	1/6/23/26	0/1/1/1
5	NAG	R	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	R	2	5	-	3/6/23/26	0/1/1/1
5	NAG	S	1	3,5	-	3/6/23/26	0/1/1/1
5	NAG	S	2	5	-	2/6/23/26	0/1/1/1
5	NAG	T	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	T	2	5	-	0/6/23/26	0/1/1/1
6	NAG	U	1	3,6	-	0/6/23/26	0/1/1/1
6	NAG	U	2	6	-	2/6/23/26	0/1/1/1
6	BMA	U	3	6	-	0/2/19/22	0/1/1/1
7	NAG	V	1	3,7	-	0/6/23/26	0/1/1/1
7	NAG	V	2	7	-	0/6/23/26	0/1/1/1
7	BMA	V	3	7	-	0/2/19/22	0/1/1/1
7	MAN	V	4	7	-	0/2/19/22	0/1/1/1
7	MAN	V	5	7	-	0/2/19/22	0/1/1/1
10	NAG	W	1	3,10	-	0/6/23/26	0/1/1/1
10	NAG	W	2	10	-	0/6/23/26	0/1/1/1
10	BMA	W	3	10	-	0/2/19/22	0/1/1/1
10	MAN	W	4	10	-	1/2/19/22	0/1/1/1
10	MAN	W	5	10	-	0/2/19/22	0/1/1/1
11	NAG	X	1	11,3	-	0/6/23/26	0/1/1/1
11	NAG	X	2	11	-	2/6/23/26	0/1/1/1
11	BMA	X	3	11	-	0/2/19/22	0/1/1/1
11	MAN	X	4	11	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	MAN	X	5	11	-	0/2/19/22	0/1/1/1
11	MAN	X	6	11	-	0/2/19/22	0/1/1/1
5	NAG	Y	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	Y	2	5	-	3/6/23/26	0/1/1/1
5	NAG	Z	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	Z	2	5	-	1/6/23/26	0/1/1/1
5	NAG	a	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	a	2	5	-	0/6/23/26	0/1/1/1
5	NAG	h	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	h	2	5	-	0/6/23/26	0/1/1/1
5	NAG	i	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	i	2	5	-	3/6/23/26	0/1/1/1
5	NAG	j	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	j	2	5	-	1/6/23/26	0/1/1/1
5	NAG	k	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	k	2	5	-	3/6/23/26	0/1/1/1
5	NAG	l	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	l	2	5	-	3/6/23/26	0/1/1/1
5	NAG	m	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	m	2	5	-	0/6/23/26	0/1/1/1
6	NAG	n	1	3,6	-	0/6/23/26	0/1/1/1
6	NAG	n	2	6	-	2/6/23/26	0/1/1/1
6	BMA	n	3	6	-	0/2/19/22	0/1/1/1
7	NAG	o	1	3,7	-	0/6/23/26	0/1/1/1
7	NAG	o	2	7	-	0/6/23/26	0/1/1/1
7	BMA	o	3	7	-	0/2/19/22	0/1/1/1
7	MAN	o	4	7	-	0/2/19/22	0/1/1/1
7	MAN	o	5	7	-	0/2/19/22	0/1/1/1
10	NAG	p	1	3,10	-	0/6/23/26	0/1/1/1
10	NAG	p	2	10	-	0/6/23/26	0/1/1/1
10	BMA	p	3	10	-	1/2/19/22	0/1/1/1
10	MAN	p	4	10	-	2/2/19/22	0/1/1/1
10	MAN	p	5	10	-	0/2/19/22	0/1/1/1
9	NAG	q	1	3,9	-	0/6/23/26	0/1/1/1
9	NAG	q	2	9	-	1/6/23/26	0/1/1/1
9	BMA	q	3	9	-	0/2/19/22	0/1/1/1
9	MAN	q	4	9	-	0/2/19/22	0/1/1/1
9	MAN	q	5	9	-	0/2/19/22	0/1/1/1
9	MAN	q	6	9	-	2/2/19/22	0/1/1/1
9	MAN	q	7	9	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MAN	q	8	9	-	1/2/19/22	0/1/1/1
5	NAG	r	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	r	2	5	-	2/6/23/26	0/1/1/1
5	NAG	s	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	s	2	5	-	1/6/23/26	0/1/1/1
5	NAG	t	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	t	2	5	-	0/6/23/26	0/1/1/1
5	NAG	u	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	u	2	5	-	1/6/23/26	0/1/1/1
5	NAG	v	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	v	2	5	-	2/6/23/26	0/1/1/1
5	NAG	w	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	w	2	5	-	1/6/23/26	0/1/1/1
5	NAG	x	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	x	2	5	-	3/6/23/26	0/1/1/1
5	NAG	y	1	3,5	-	3/6/23/26	0/1/1/1
5	NAG	y	2	5	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	U	3	BMA	C1-O5-C5	6.53	120.94	112.19
6	n	3	BMA	C1-O5-C5	6.52	120.93	112.19
11	X	3	BMA	C1-O5-C5	6.52	120.93	112.19
6	B	3	BMA	C1-O5-C5	6.47	120.86	112.19
5	l	2	NAG	C1-C2-N2	3.82	116.45	110.43
5	l	2	NAG	C2-N2-C7	3.63	127.77	122.90
5	a	2	NAG	C1-O5-C5	3.60	117.01	112.19
11	X	3	BMA	C3-C4-C5	3.52	116.62	110.23
5	i	2	NAG	C1-O5-C5	3.46	116.82	112.19
5	j	2	NAG	C1-O5-C5	3.44	116.79	112.19
11	X	4	MAN	C1-O5-C5	3.31	116.63	112.19
9	q	2	NAG	C2-N2-C7	3.26	127.27	122.90
6	n	3	BMA	C3-C4-C5	2.84	115.39	110.23
5	t	2	NAG	C1-C2-N2	2.82	114.87	110.43
6	U	3	BMA	C3-C4-C5	2.82	115.34	110.23
5	K	2	NAG	C1-C2-N2	2.80	114.85	110.43
6	B	3	BMA	C3-C4-C5	2.79	115.29	110.23
11	X	3	BMA	O4-C4-C3	-2.67	104.08	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	X	3	BMA	C2-C3-C4	2.63	115.49	110.86
5	w	2	NAG	C1-C2-N2	2.62	114.57	110.43
5	Q	2	NAG	C1-C2-N2	2.61	114.55	110.43
5	G	2	NAG	C1-C2-N2	2.51	114.39	110.43
5	j	1	NAG	O5-C1-C2	-2.46	107.49	111.29
5	i	2	NAG	C2-N2-C7	2.42	126.14	122.90
11	X	2	NAG	C4-C3-C2	-2.39	107.52	111.02
5	v	2	NAG	C1-C2-N2	2.34	114.12	110.43
10	p	4	MAN	C1-O5-C5	2.32	115.30	112.19
11	X	5	MAN	C1-O5-C5	2.32	115.29	112.19
5	m	1	NAG	O5-C1-C2	-2.30	107.72	111.29
5	P	2	NAG	C1-C2-N2	2.30	114.06	110.43
5	y	2	NAG	C2-N2-C7	2.27	125.94	122.90
5	k	2	NAG	C2-N2-C7	2.23	125.89	122.90
5	h	2	NAG	O5-C1-C2	-2.22	107.85	111.29
5	S	2	NAG	C2-N2-C7	2.22	125.88	122.90
5	A	1	NAG	O5-C1-C2	-2.22	107.86	111.29
6	n	3	BMA	C2-C3-C4	2.18	114.70	110.86
9	E	6	MAN	O3-C3-C2	-2.17	105.62	110.05
6	U	3	BMA	C2-C3-C4	2.17	114.68	110.86
5	h	1	NAG	O5-C1-C2	-2.17	107.94	111.29
5	T	1	NAG	O5-C1-C2	-2.17	107.94	111.29
6	B	3	BMA	C2-C3-C4	2.16	114.66	110.86
5	Z	1	NAG	O5-C1-C2	-2.15	107.96	111.29
6	n	3	BMA	O4-C4-C3	-2.12	105.38	110.38
5	y	1	NAG	C2-N2-C7	2.11	125.73	122.90
6	B	3	BMA	O4-C4-C3	-2.11	105.39	110.38
6	U	3	BMA	O4-C4-C3	-2.11	105.40	110.38
6	n	2	NAG	O5-C1-C2	-2.10	108.04	111.29
11	X	3	BMA	O5-C5-C4	2.10	115.93	110.83
5	h	1	NAG	O4-C4-C3	-2.09	105.44	110.38
5	R	2	NAG	C2-N2-C7	2.09	125.70	122.90
11	X	6	MAN	C1-O5-C5	2.08	114.98	112.19
5	x	2	NAG	C2-N2-C7	2.08	125.68	122.90
5	i	1	NAG	O5-C1-C2	-2.08	108.08	111.29
6	U	2	NAG	O5-C1-C2	-2.07	108.08	111.29
5	K	2	NAG	C1-O5-C5	2.07	114.96	112.19
5	a	2	NAG	C1-C2-N2	2.06	113.69	110.43
5	t	2	NAG	C1-O5-C5	2.06	114.94	112.19
6	B	2	NAG	O5-C1-C2	-2.05	108.12	111.29
9	E	2	NAG	C2-N2-C7	2.05	125.65	122.90
5	Y	1	NAG	O4-C4-C3	-2.04	105.57	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	Y	2	NAG	O5-C1-C2	-2.03	108.15	111.29
5	r	2	NAG	C2-N2-C7	2.01	125.60	122.90
6	n	2	NAG	C2-N2-C7	2.01	125.60	122.90

There are no chirality outliers.

All (86) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	k	1	NAG	O7-C7-N2-C2
5	l	2	NAG	C1-C2-N2-C7
5	k	1	NAG	C8-C7-N2-C2
10	p	4	MAN	O5-C5-C6-O6
5	G	1	NAG	C8-C7-N2-C2
5	G	1	NAG	O7-C7-N2-C2
5	s	1	NAG	C8-C7-N2-C2
5	R	1	NAG	C8-C7-N2-C2
5	S	1	NAG	C8-C7-N2-C2
5	S	1	NAG	O7-C7-N2-C2
5	l	2	NAG	C8-C7-N2-C2
5	l	2	NAG	O7-C7-N2-C2
5	s	1	NAG	O7-C7-N2-C2
5	x	1	NAG	C8-C7-N2-C2
5	R	1	NAG	O7-C7-N2-C2
10	p	4	MAN	C4-C5-C6-O6
5	x	1	NAG	O7-C7-N2-C2
5	k	2	NAG	O5-C5-C6-O6
5	G	2	NAG	O5-C5-C6-O6
5	P	2	NAG	O5-C5-C6-O6
5	R	2	NAG	O5-C5-C6-O6
5	x	2	NAG	O5-C5-C6-O6
5	s	2	NAG	O5-C5-C6-O6
5	v	2	NAG	O5-C5-C6-O6
5	F	2	NAG	O5-C5-C6-O6
5	Q	2	NAG	O5-C5-C6-O6
5	S	2	NAG	O5-C5-C6-O6
5	j	2	NAG	O5-C5-C6-O6
5	r	2	NAG	O5-C5-C6-O6
5	w	2	NAG	O5-C5-C6-O6
5	i	2	NAG	O5-C5-C6-O6
5	Z	2	NAG	O5-C5-C6-O6
9	q	8	MAN	O5-C5-C6-O6
5	Y	2	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	y	1	NAG	O5-C5-C6-O6
5	l	1	NAG	O5-C5-C6-O6
5	t	1	NAG	O5-C5-C6-O6
8	D	5	MAN	O5-C5-C6-O6
10	W	4	MAN	O5-C5-C6-O6
10	p	3	BMA	O5-C5-C6-O6
5	K	1	NAG	O5-C5-C6-O6
5	S	1	NAG	O5-C5-C6-O6
5	a	1	NAG	O5-C5-C6-O6
5	S	2	NAG	C3-C2-N2-C7
9	q	6	MAN	C4-C5-C6-O6
5	Y	2	NAG	C1-C2-N2-C7
5	h	1	NAG	C1-C2-N2-C7
6	B	2	NAG	C1-C2-N2-C7
6	U	2	NAG	C1-C2-N2-C7
6	n	2	NAG	C1-C2-N2-C7
11	X	2	NAG	C1-C2-N2-C7
9	q	6	MAN	O5-C5-C6-O6
5	F	1	NAG	C3-C2-N2-C7
5	F	2	NAG	C3-C2-N2-C7
5	O	1	NAG	C3-C2-N2-C7
5	P	2	NAG	C3-C2-N2-C7
5	R	2	NAG	C3-C2-N2-C7
5	Y	1	NAG	C3-C2-N2-C7
5	Y	2	NAG	C3-C2-N2-C7
5	h	1	NAG	C3-C2-N2-C7
5	i	2	NAG	C3-C2-N2-C7
5	k	2	NAG	C3-C2-N2-C7
5	r	1	NAG	C3-C2-N2-C7
5	r	2	NAG	C3-C2-N2-C7
5	u	1	NAG	C3-C2-N2-C7
5	v	2	NAG	C3-C2-N2-C7
5	x	2	NAG	C3-C2-N2-C7
5	y	1	NAG	C3-C2-N2-C7
5	y	2	NAG	C3-C2-N2-C7
9	E	2	NAG	C3-C2-N2-C7
9	q	2	NAG	C3-C2-N2-C7
11	X	2	NAG	C3-C2-N2-C7
5	O	1	NAG	C1-C2-N2-C7
5	R	2	NAG	C1-C2-N2-C7
5	Y	1	NAG	C1-C2-N2-C7
5	i	2	NAG	C1-C2-N2-C7

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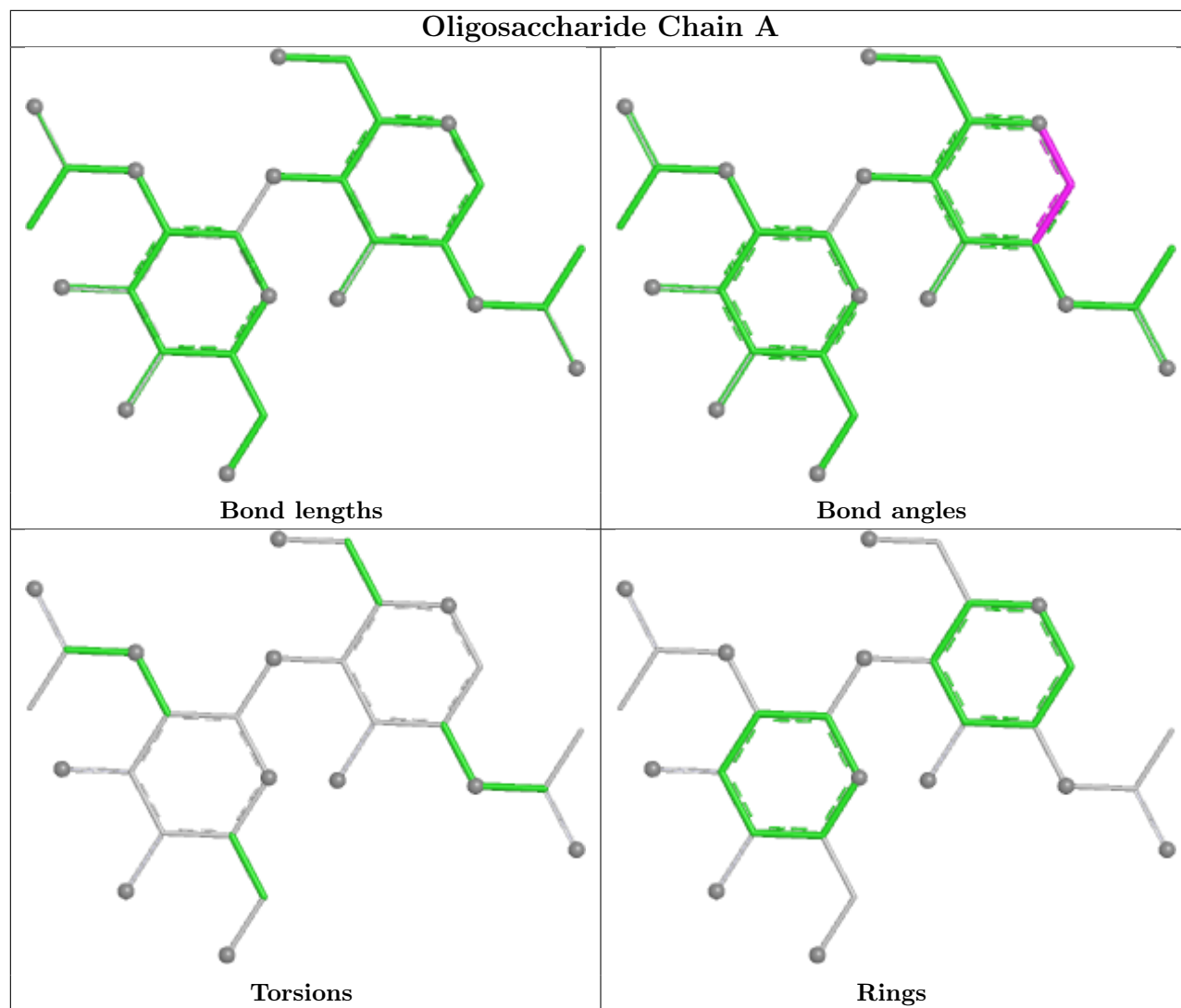
Mol	Chain	Res	Type	Atoms
5	k	2	NAG	C1-C2-N2-C7
5	u	1	NAG	C1-C2-N2-C7
5	x	2	NAG	C1-C2-N2-C7
5	y	1	NAG	C1-C2-N2-C7
9	E	2	NAG	C1-C2-N2-C7
5	O	2	NAG	C3-C2-N2-C7
5	u	2	NAG	C3-C2-N2-C7
6	B	2	NAG	C3-C2-N2-C7
6	U	2	NAG	C3-C2-N2-C7
6	n	2	NAG	C3-C2-N2-C7

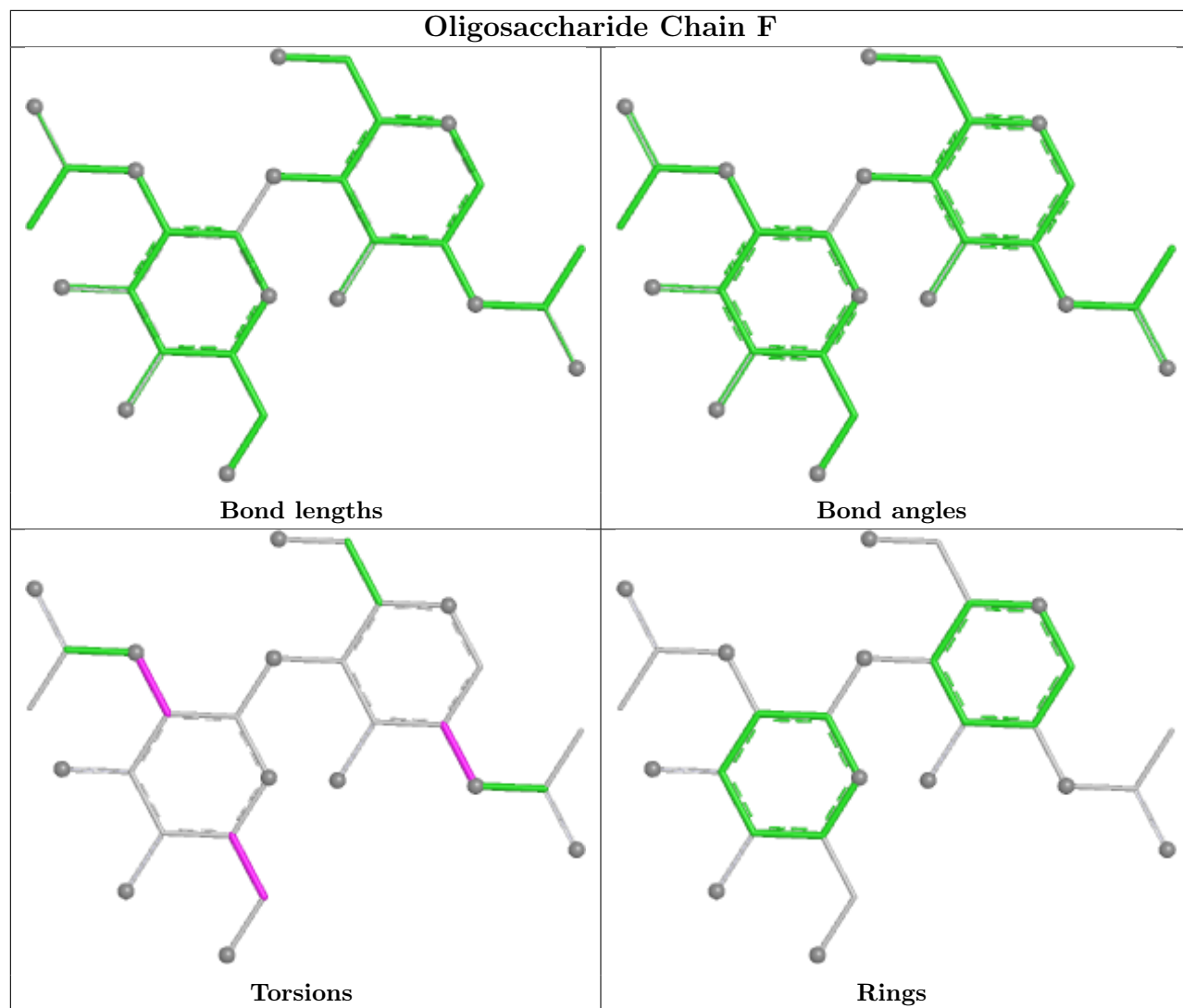
There are no ring outliers.

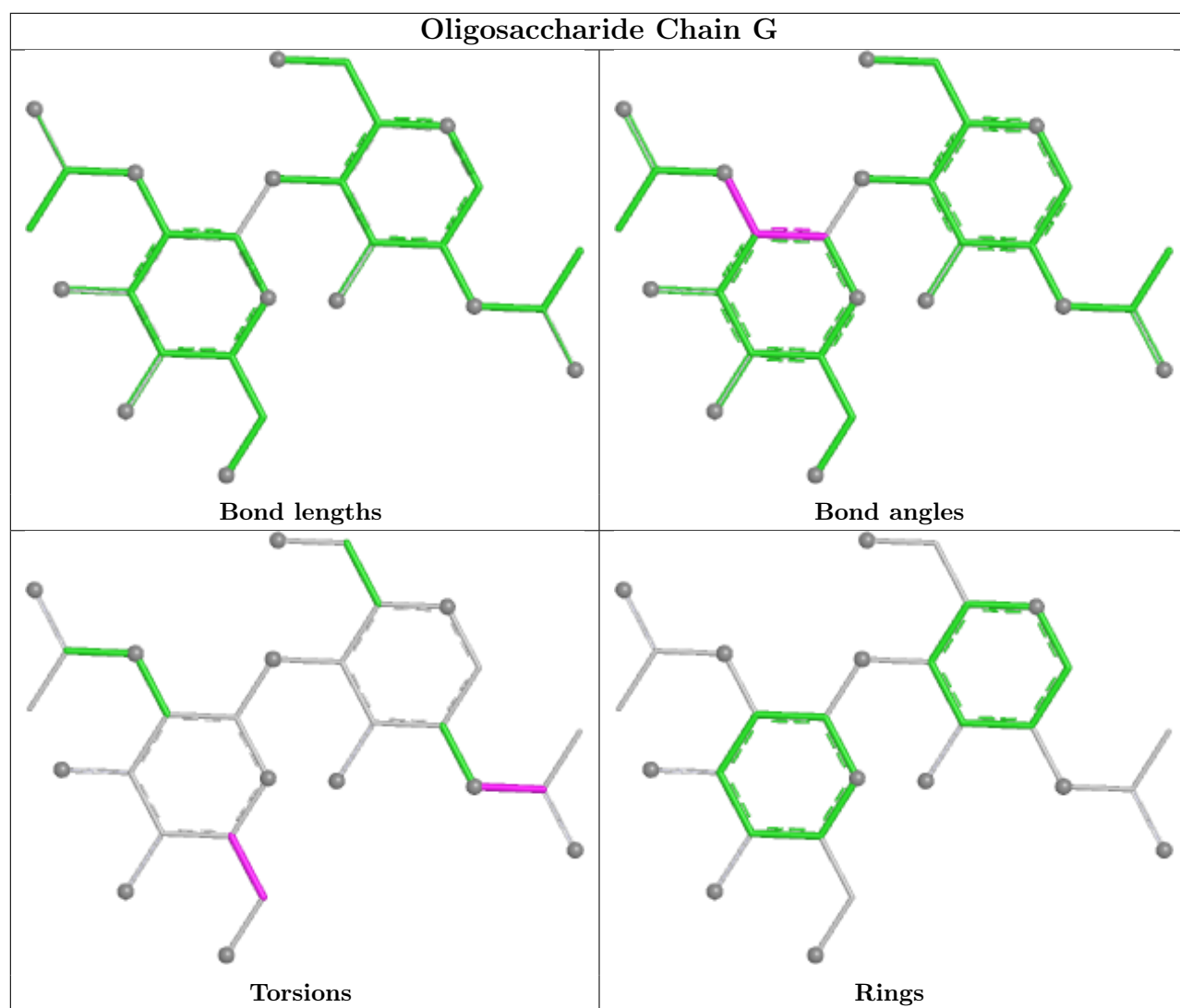
22 monomers are involved in 25 short contacts:

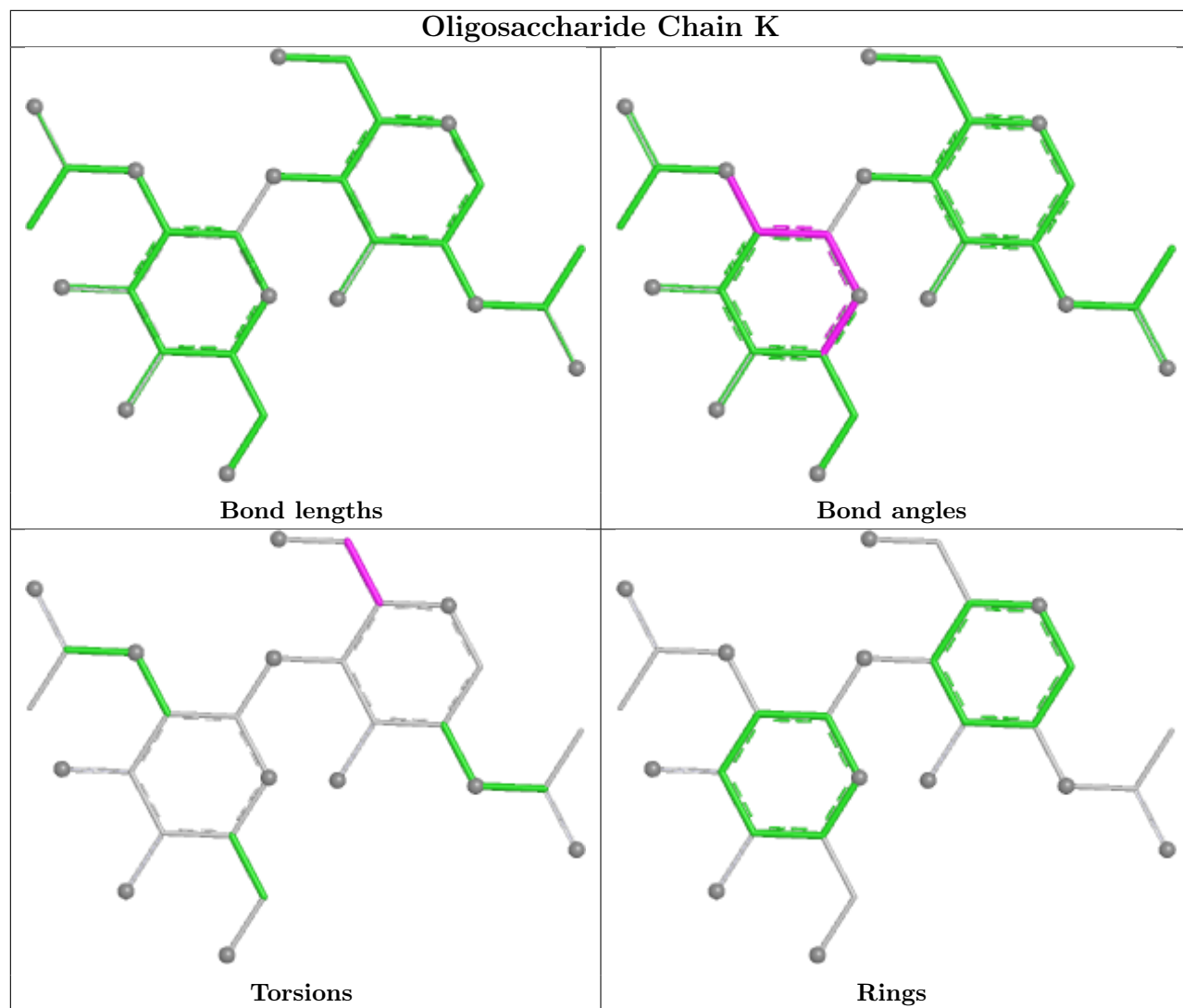
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	v	2	NAG	1	0
9	q	2	NAG	2	0
10	W	5	MAN	1	0
5	t	1	NAG	3	0
5	P	1	NAG	1	0
5	K	2	NAG	2	0
10	p	1	NAG	2	0
10	p	4	MAN	1	0
8	D	5	MAN	2	0
10	p	3	BMA	1	0
11	X	5	MAN	1	0
5	y	1	NAG	1	0
5	v	1	NAG	1	0
11	X	4	MAN	1	0
10	p	5	MAN	3	0
5	G	1	NAG	1	0
5	P	2	NAG	1	0
10	W	1	NAG	1	0
5	K	1	NAG	2	0
9	E	2	NAG	2	0
8	D	1	NAG	1	0
5	t	2	NAG	2	0

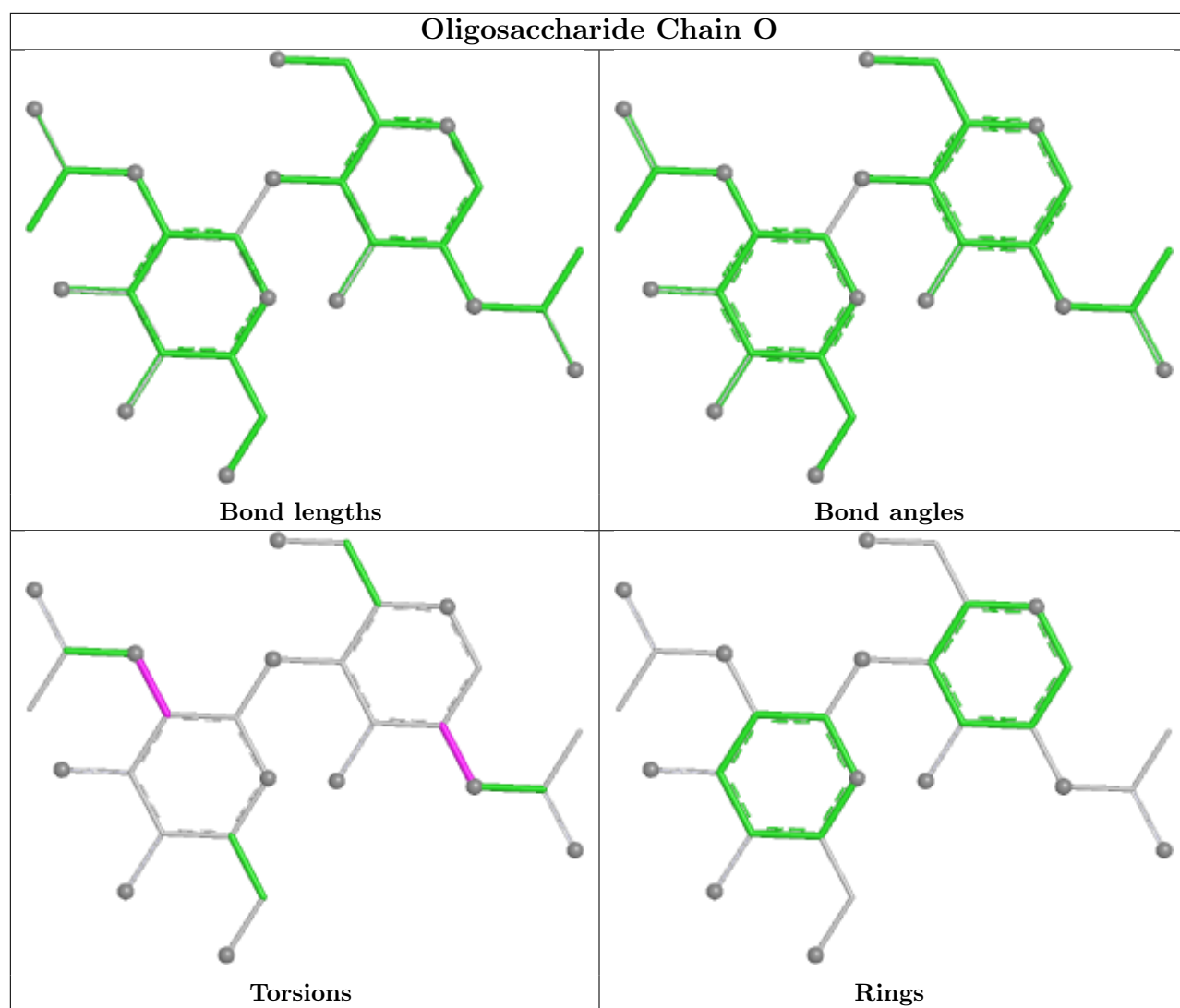
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

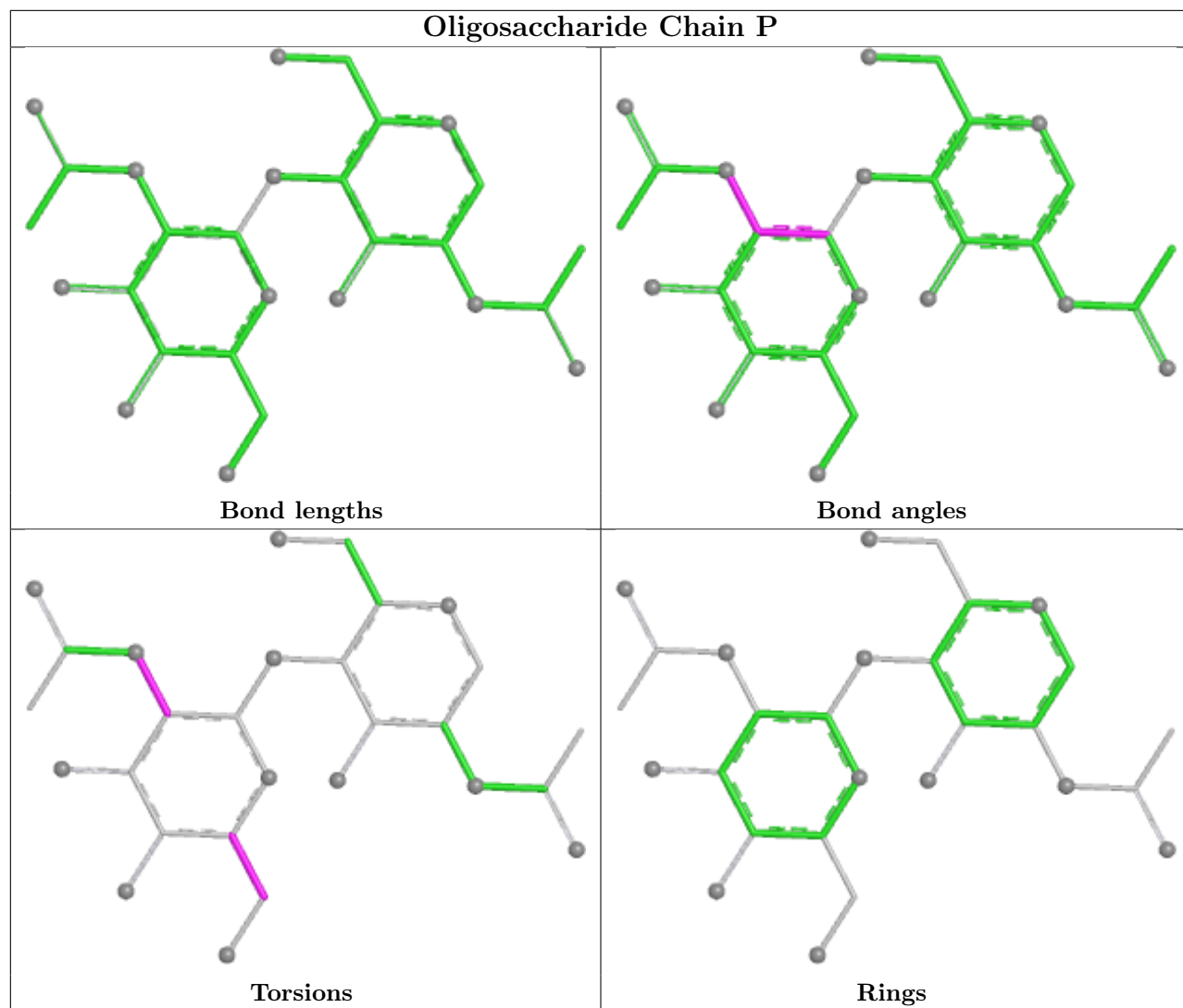


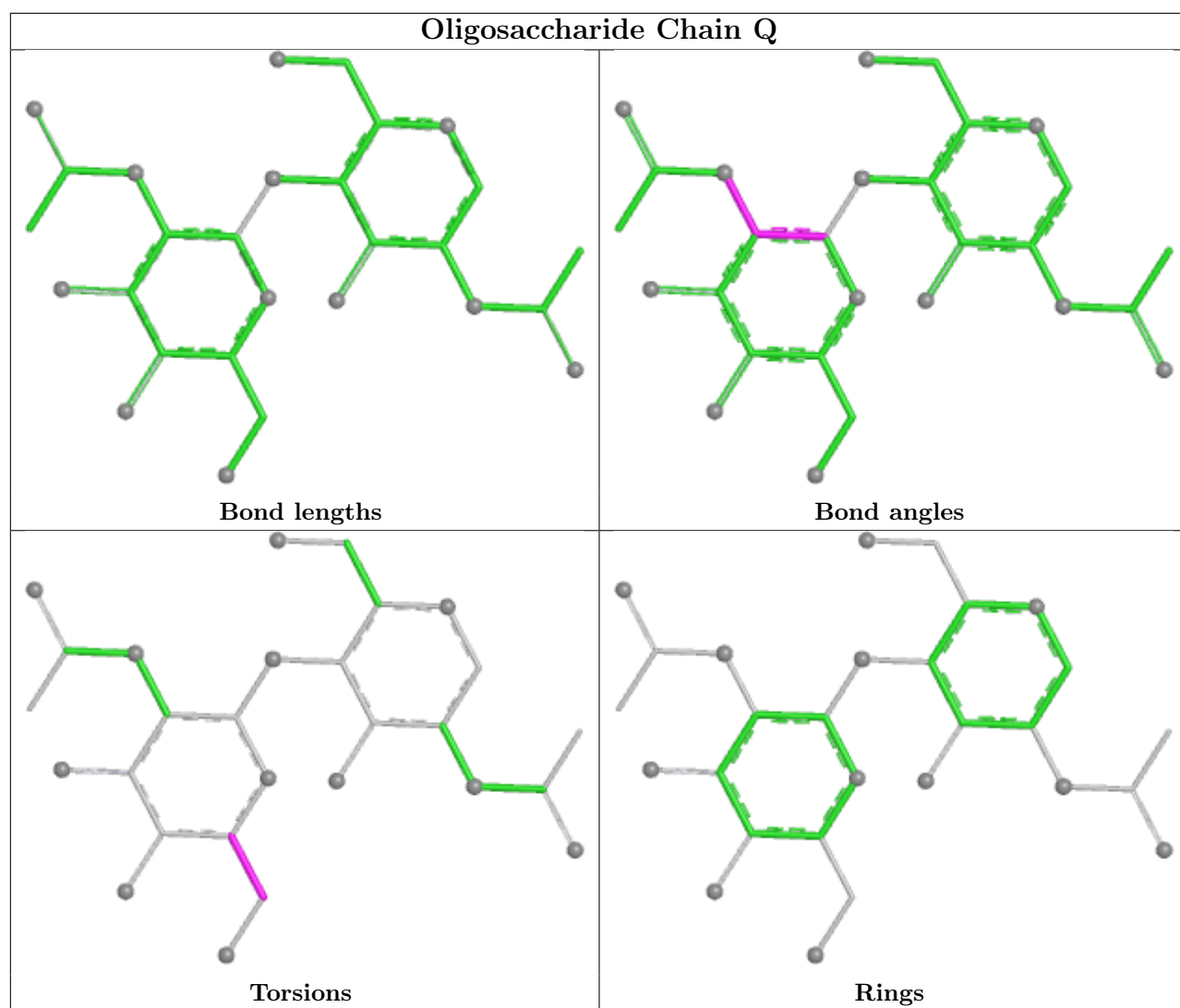


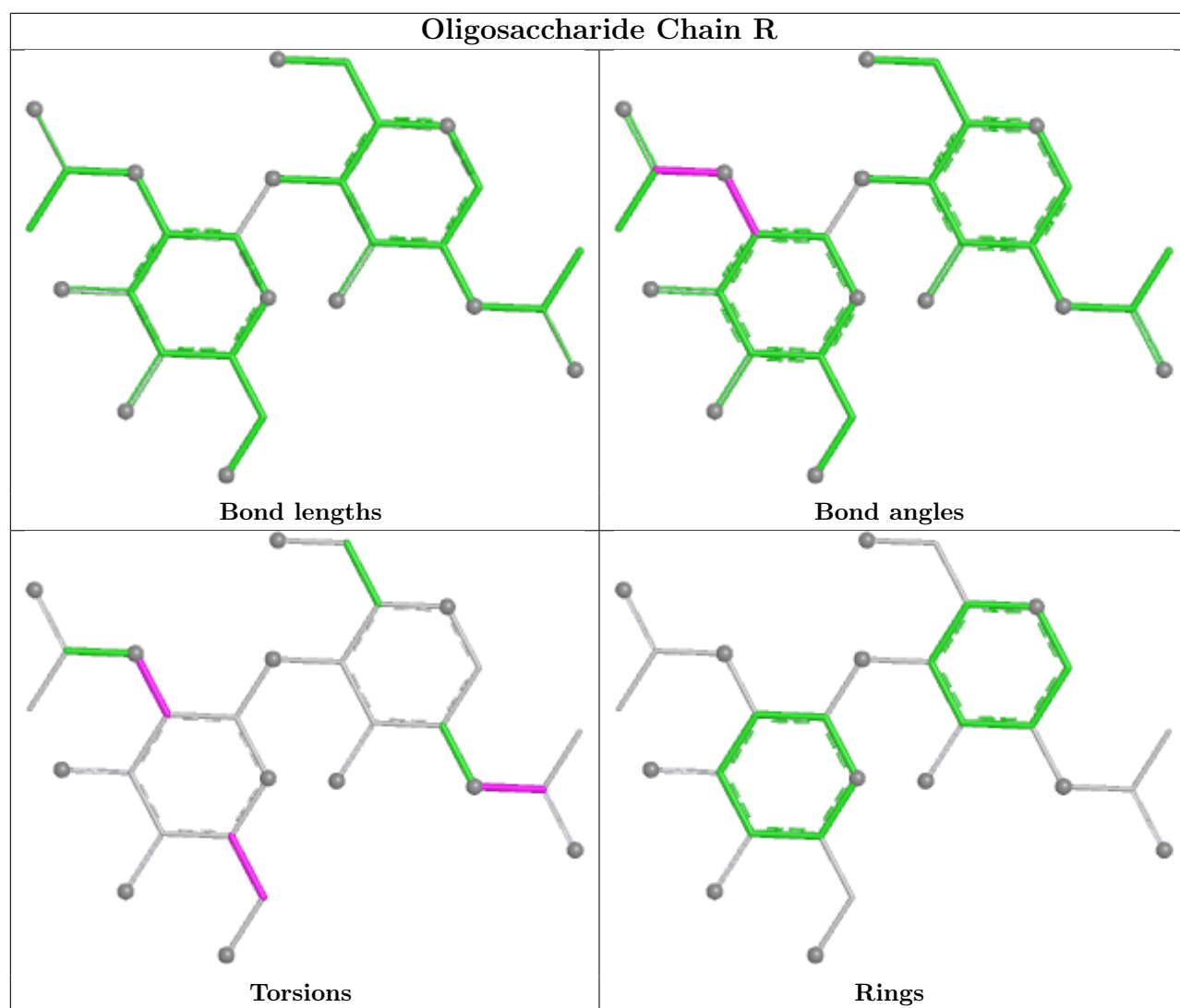


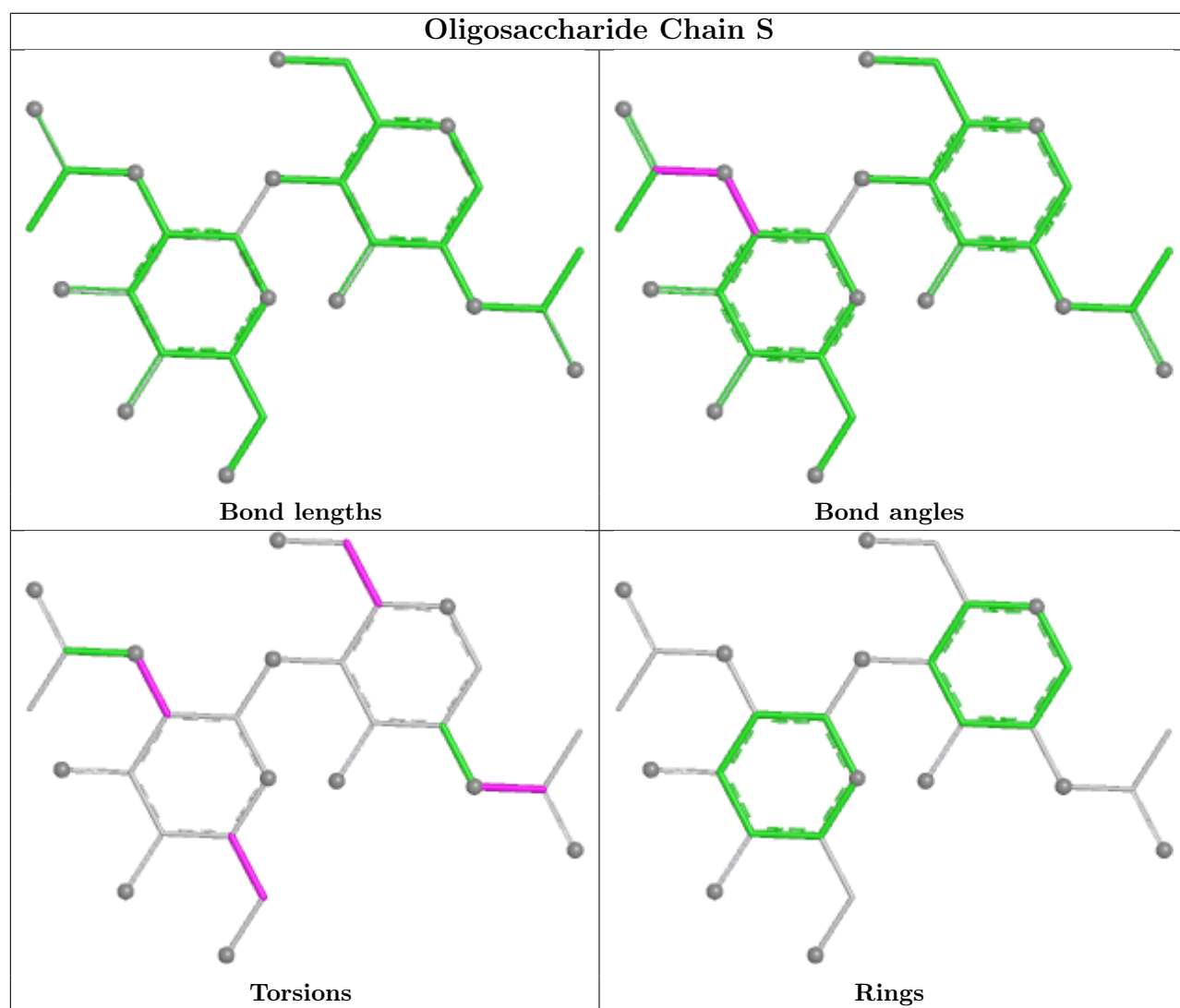


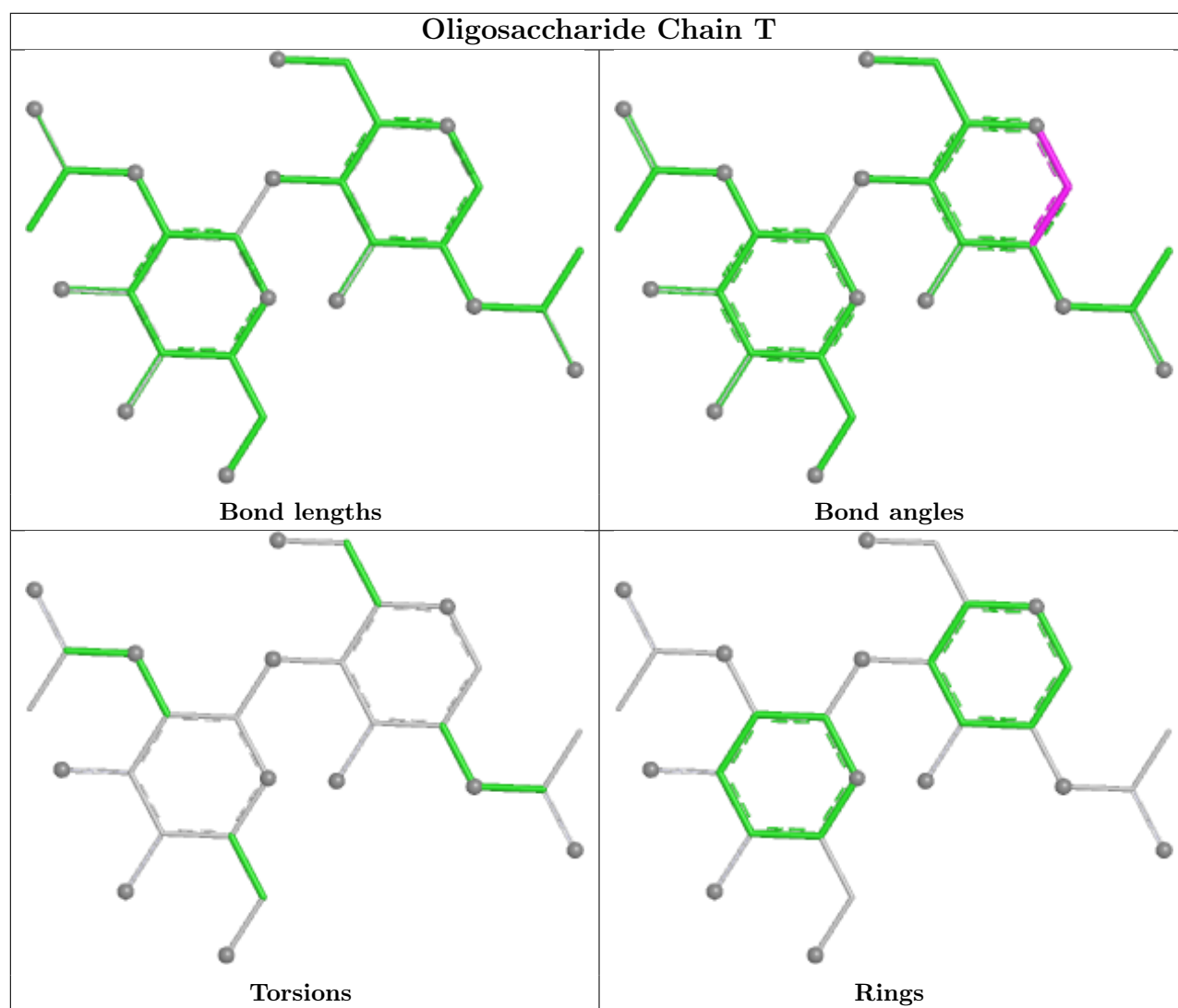


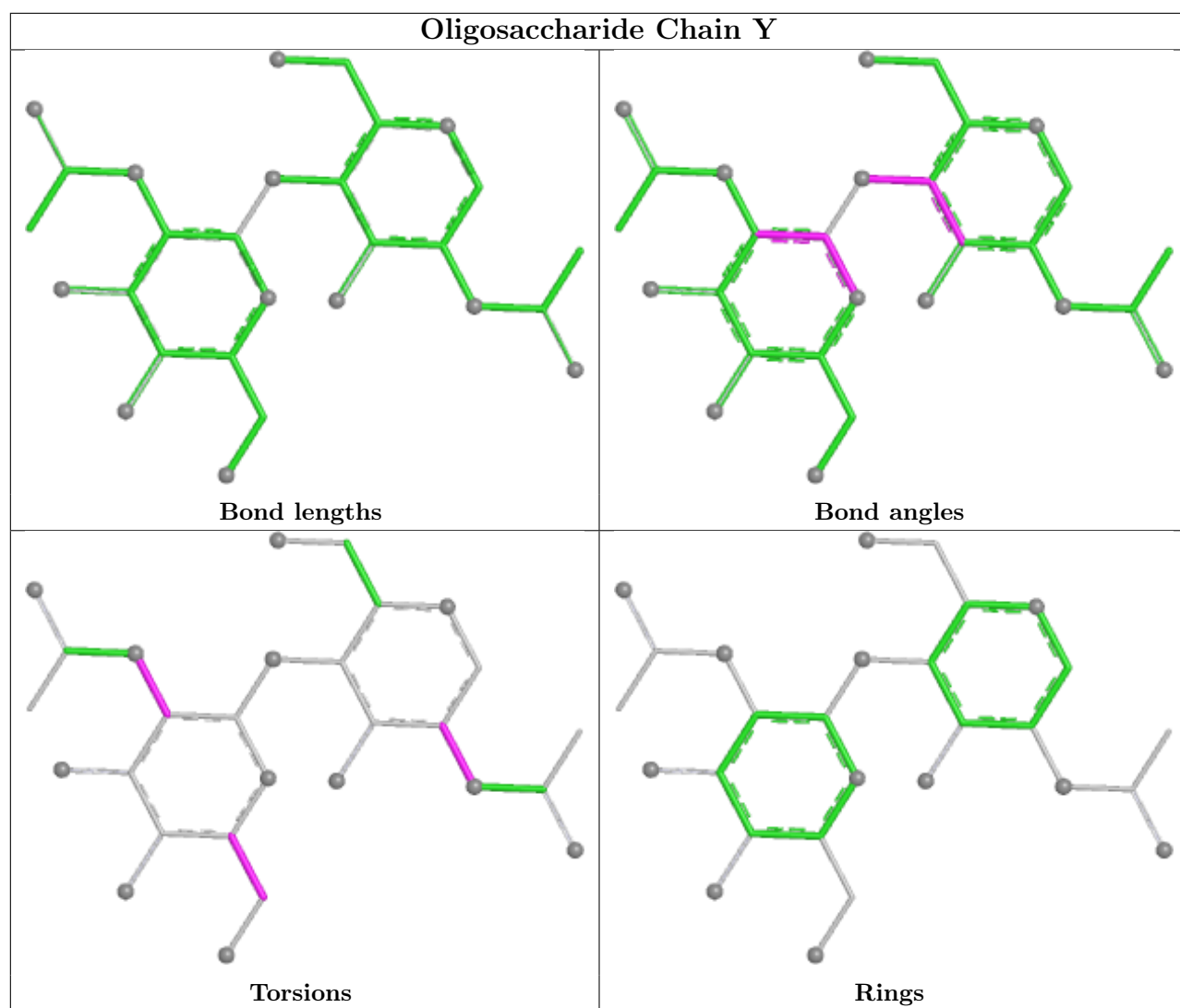


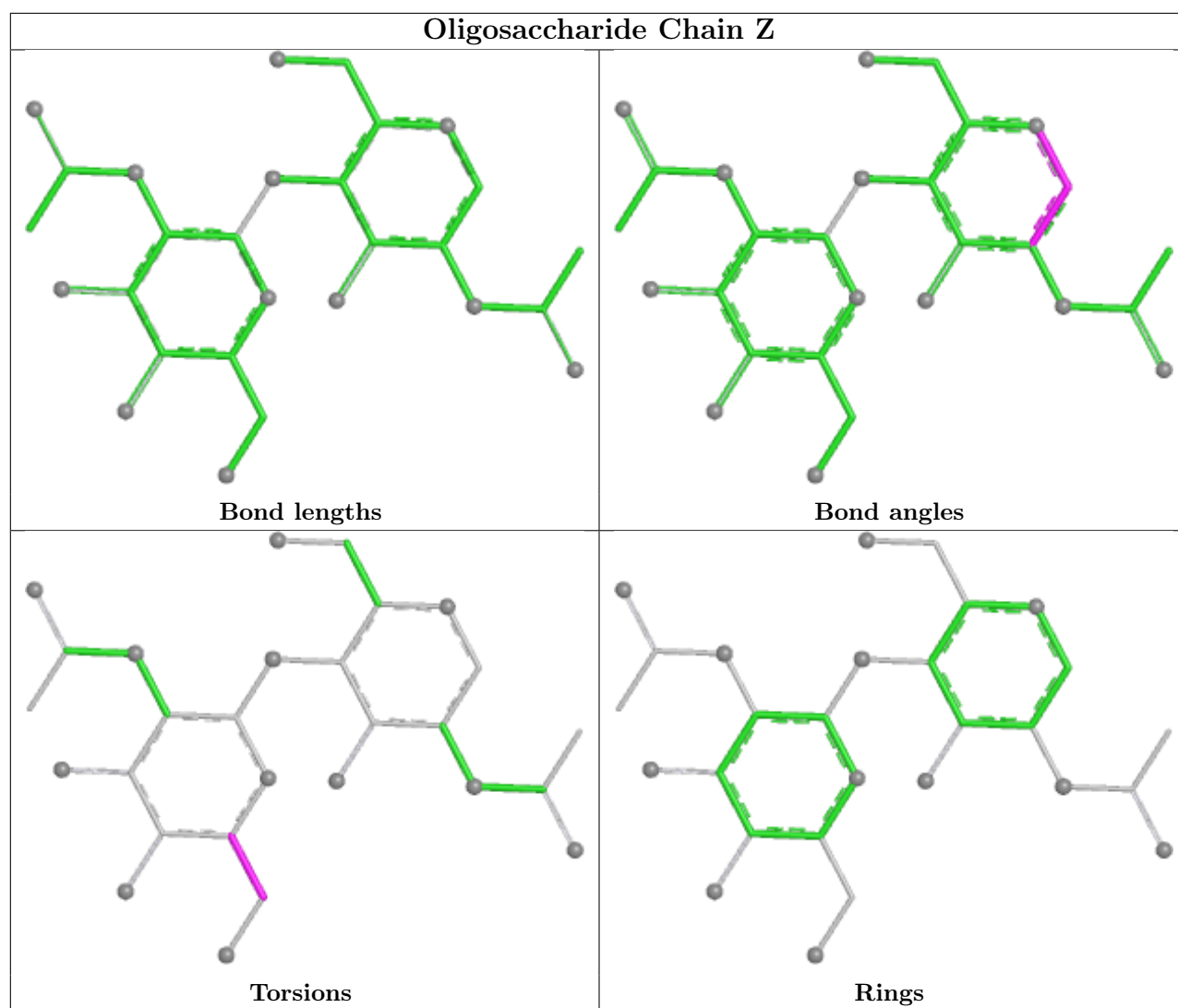


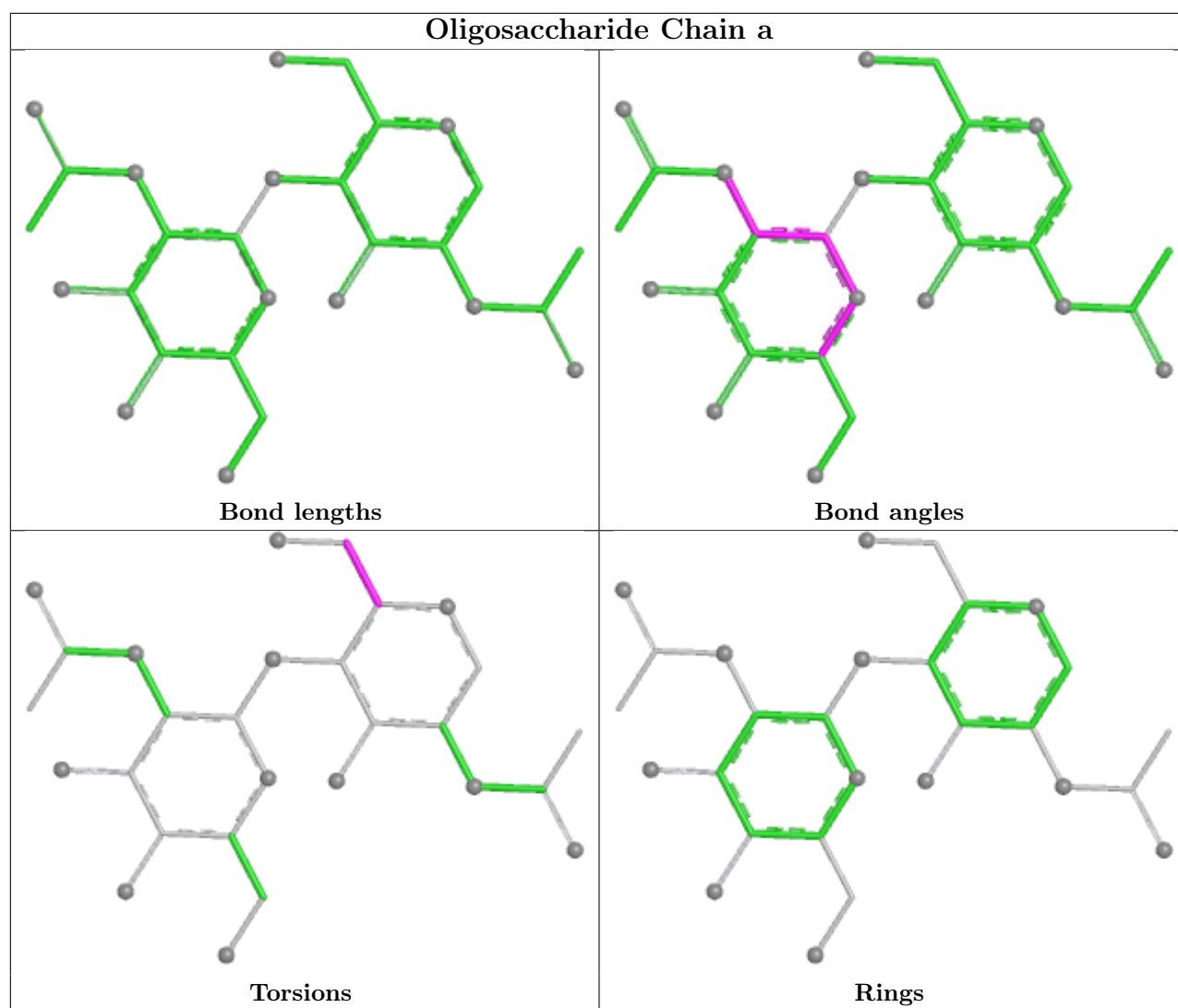


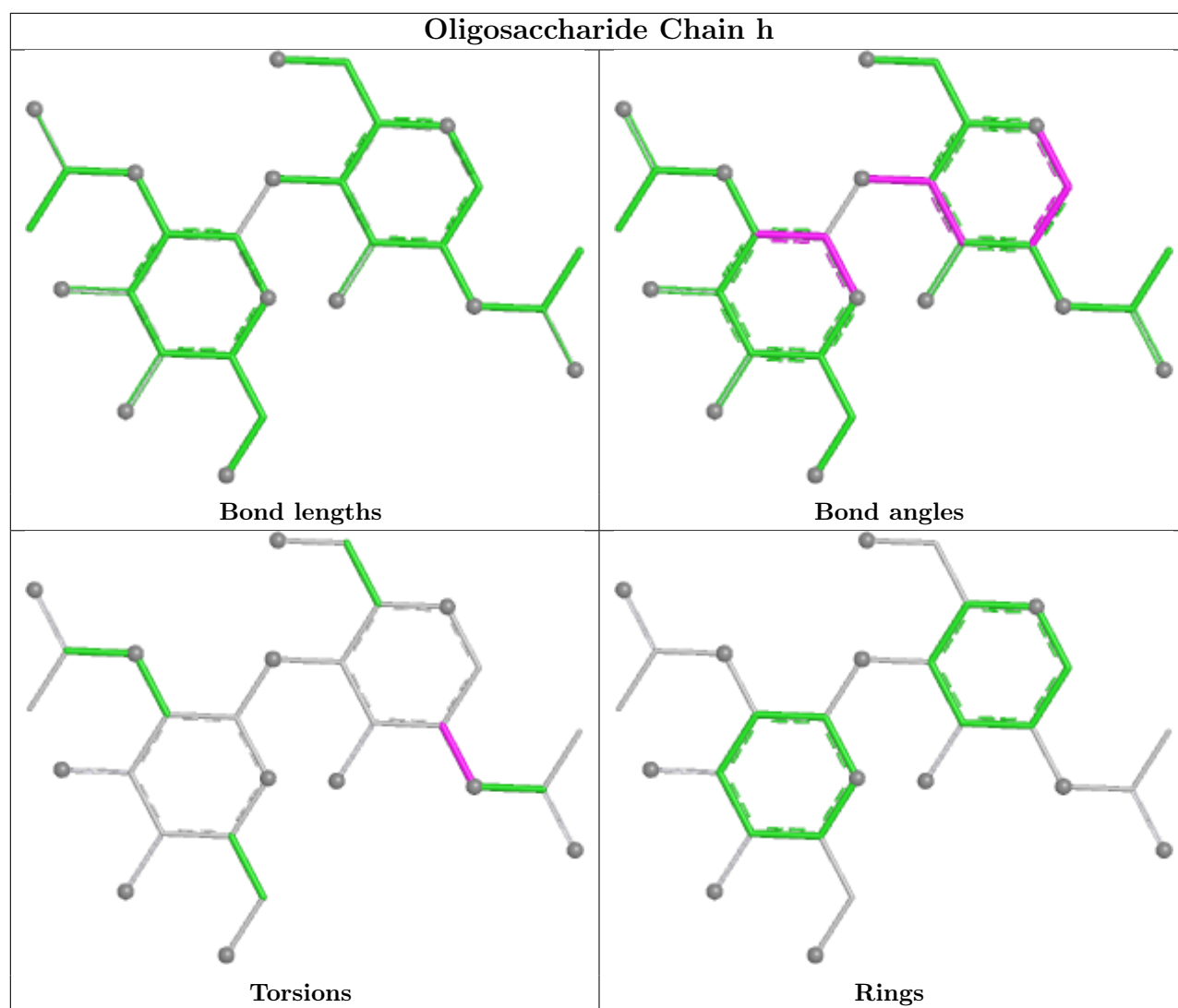


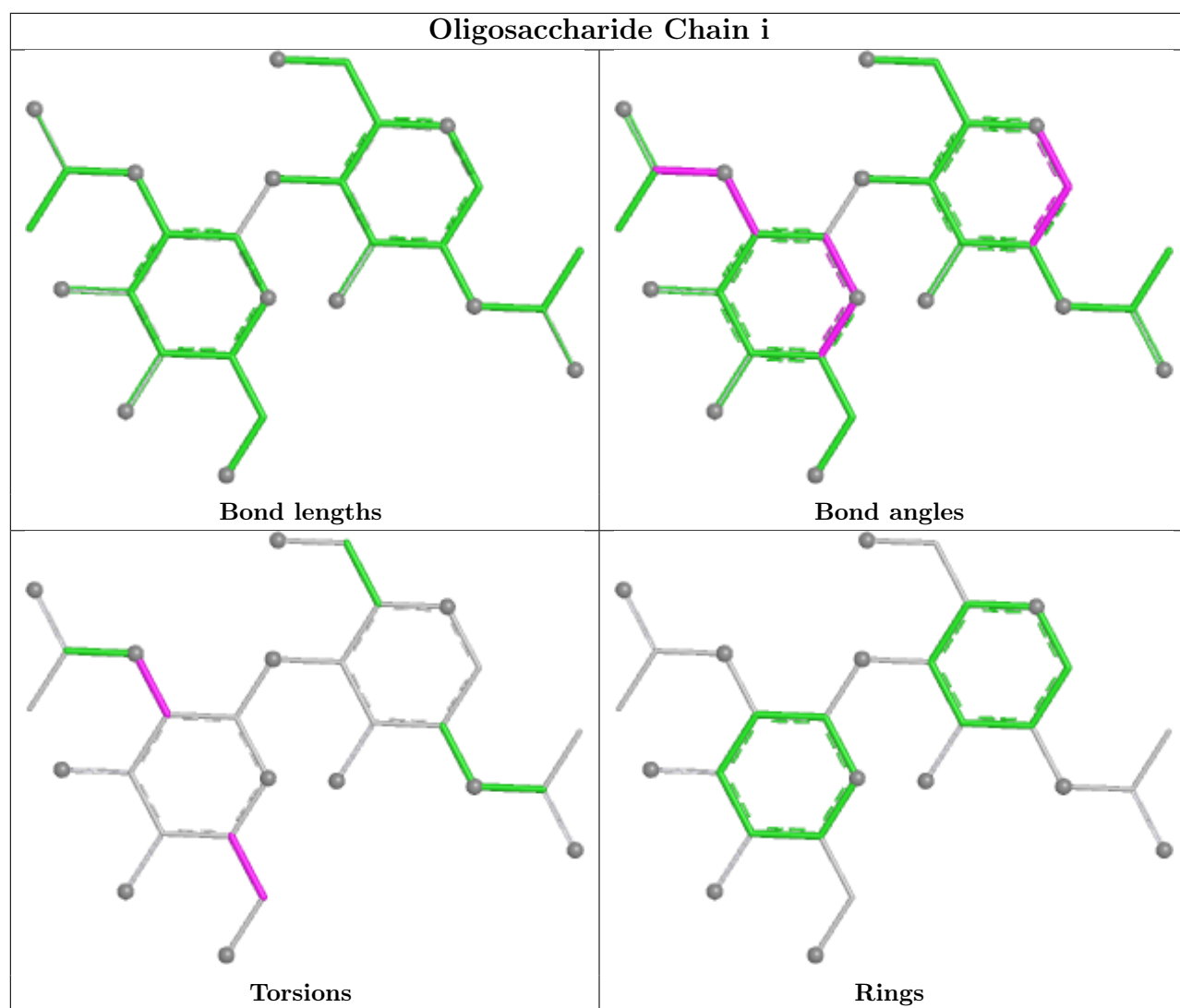


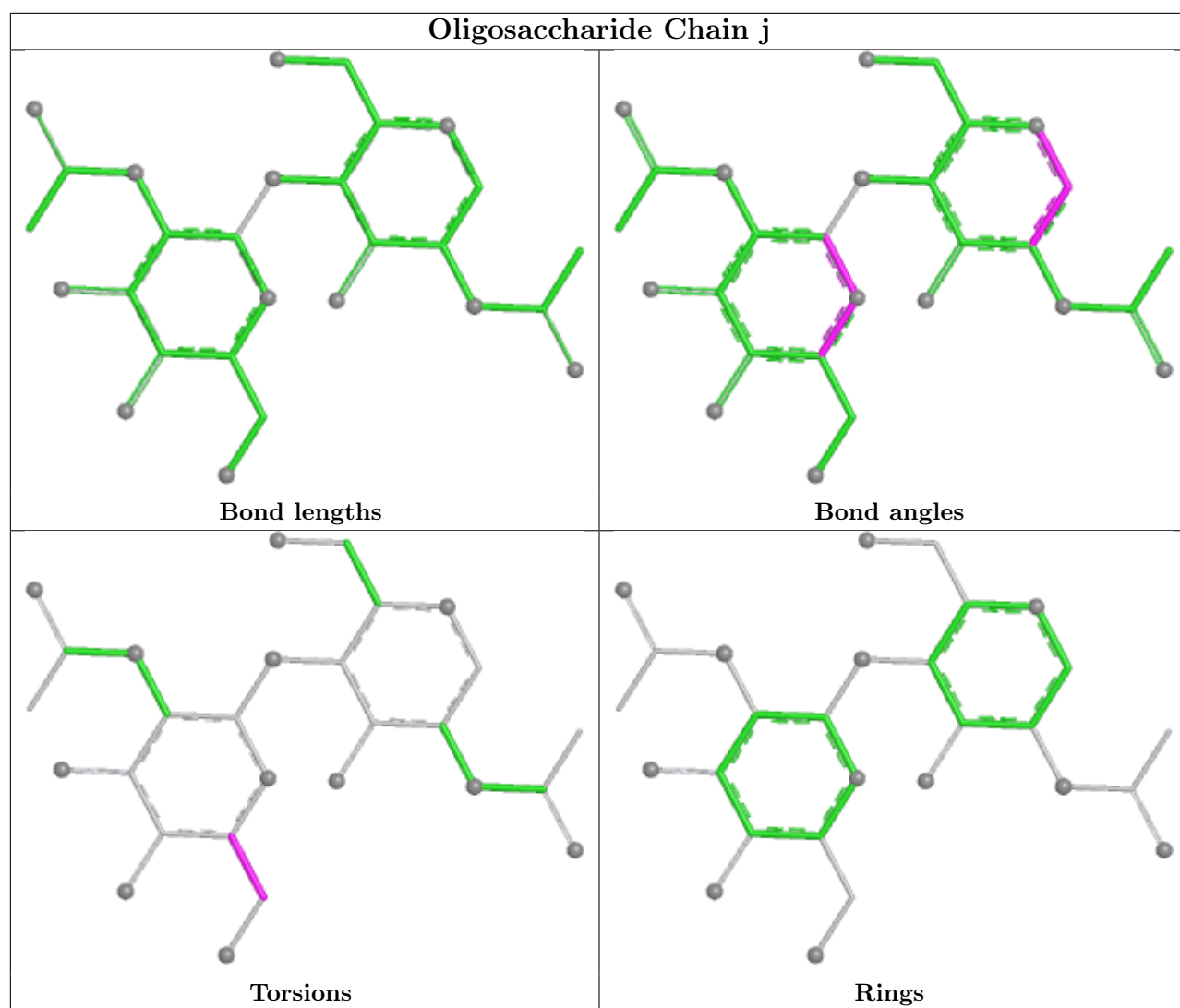


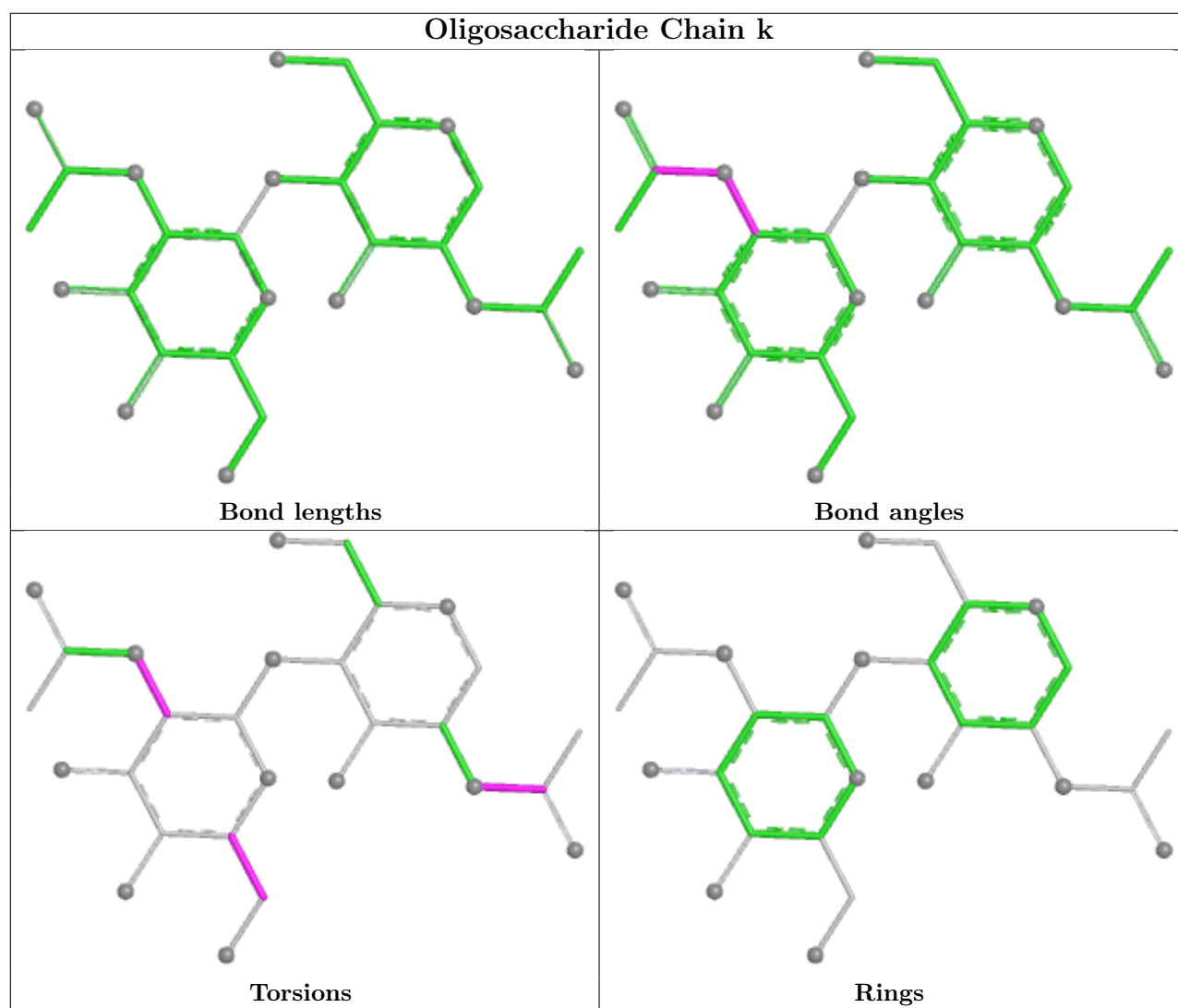


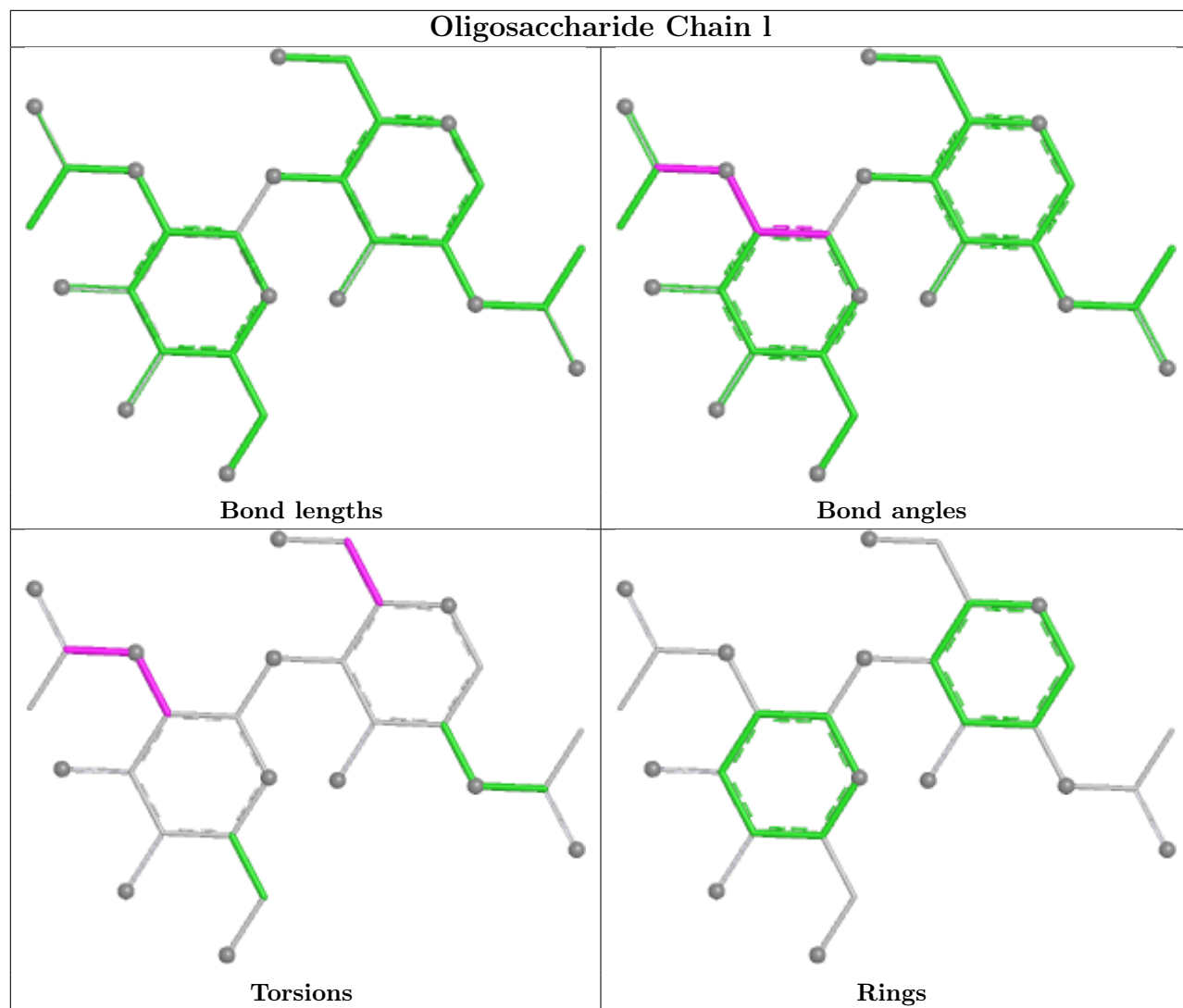


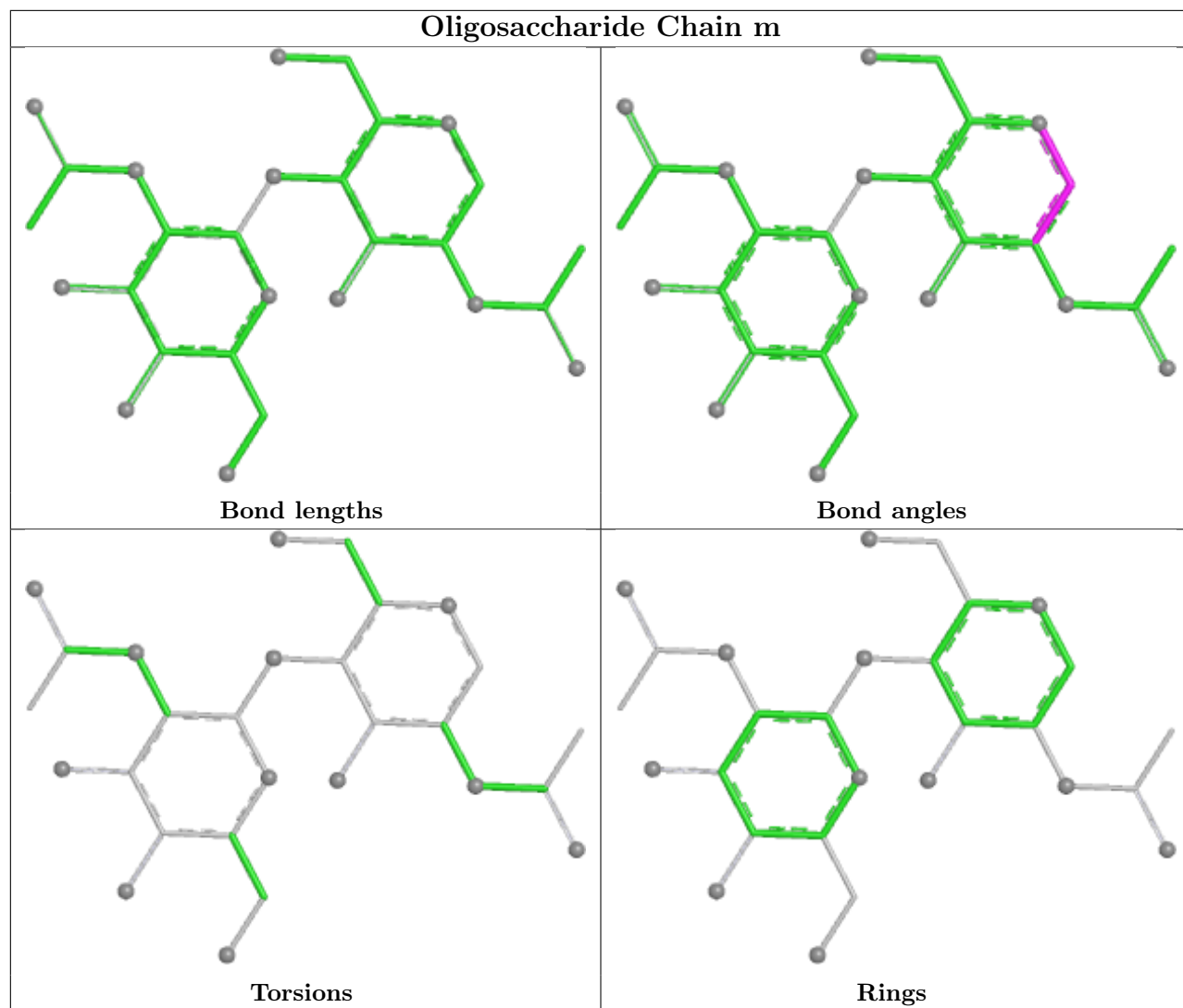


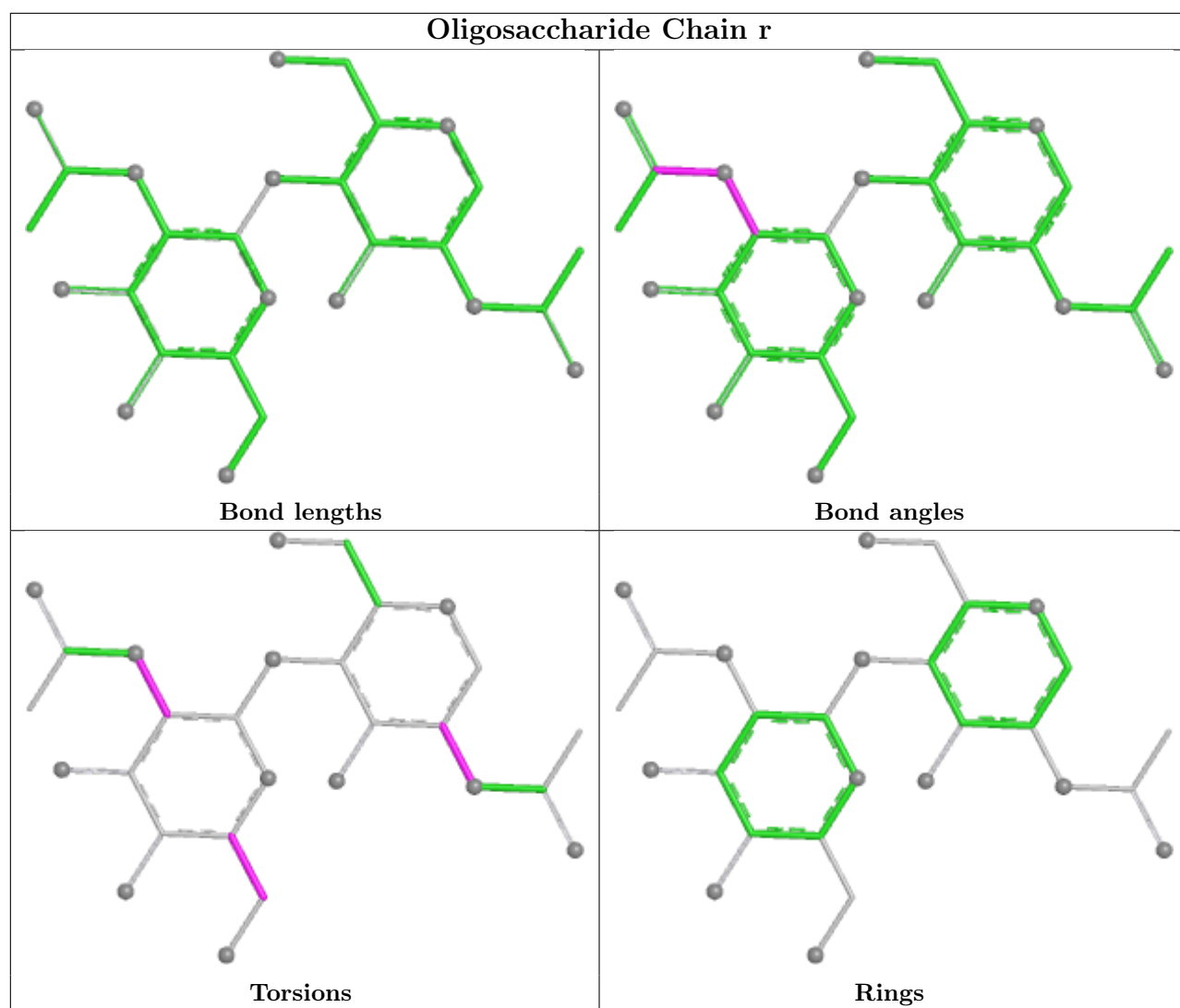


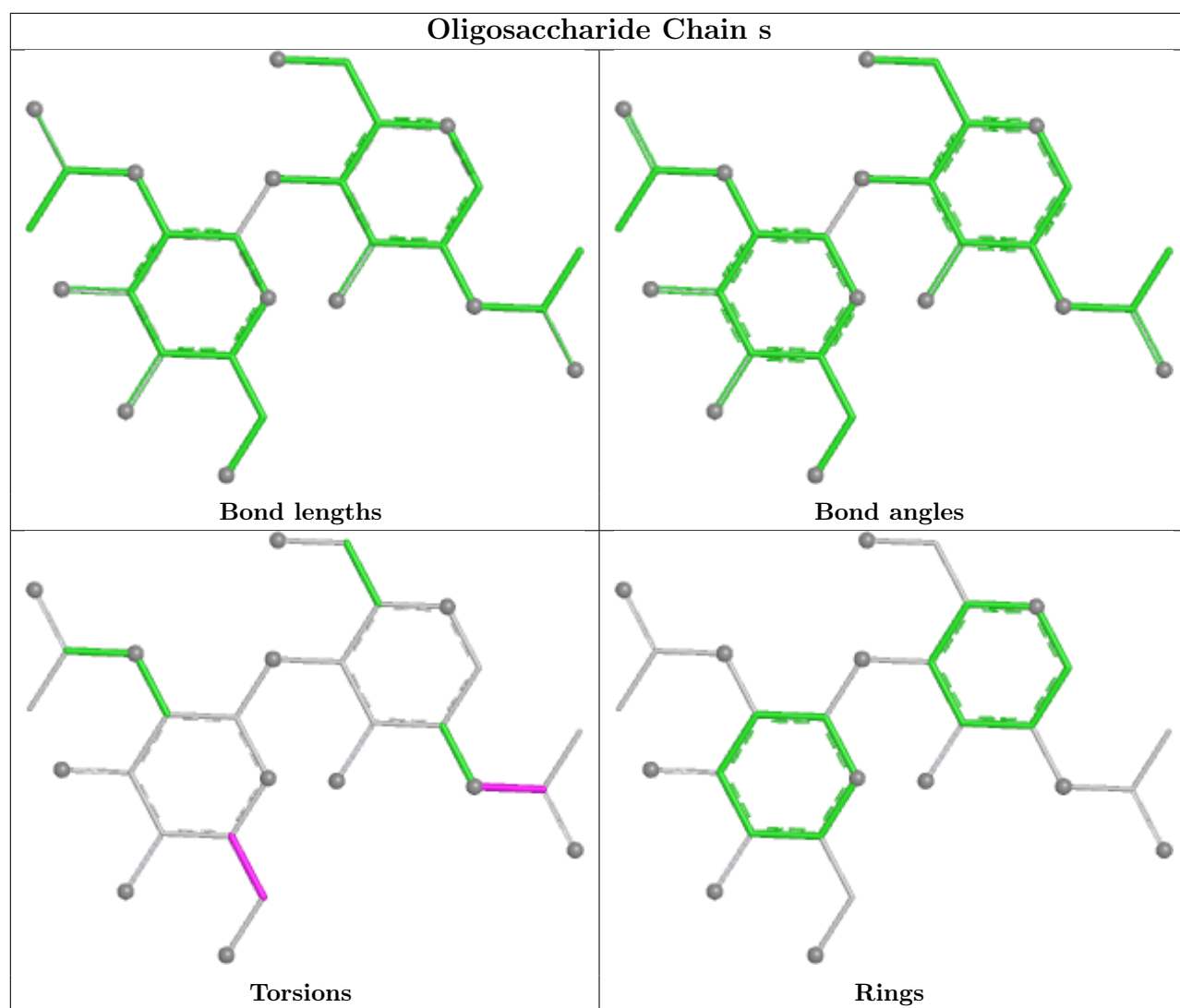


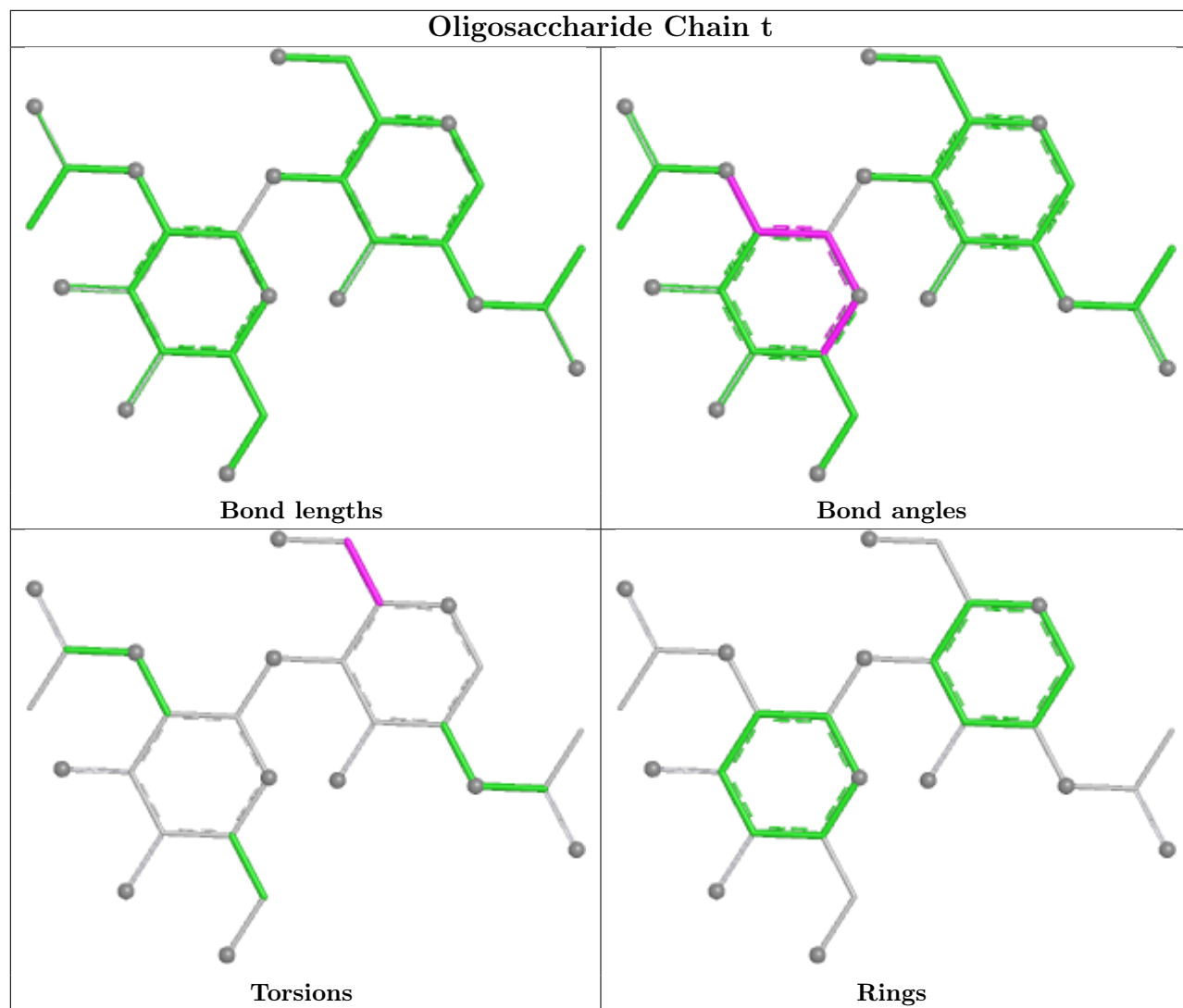


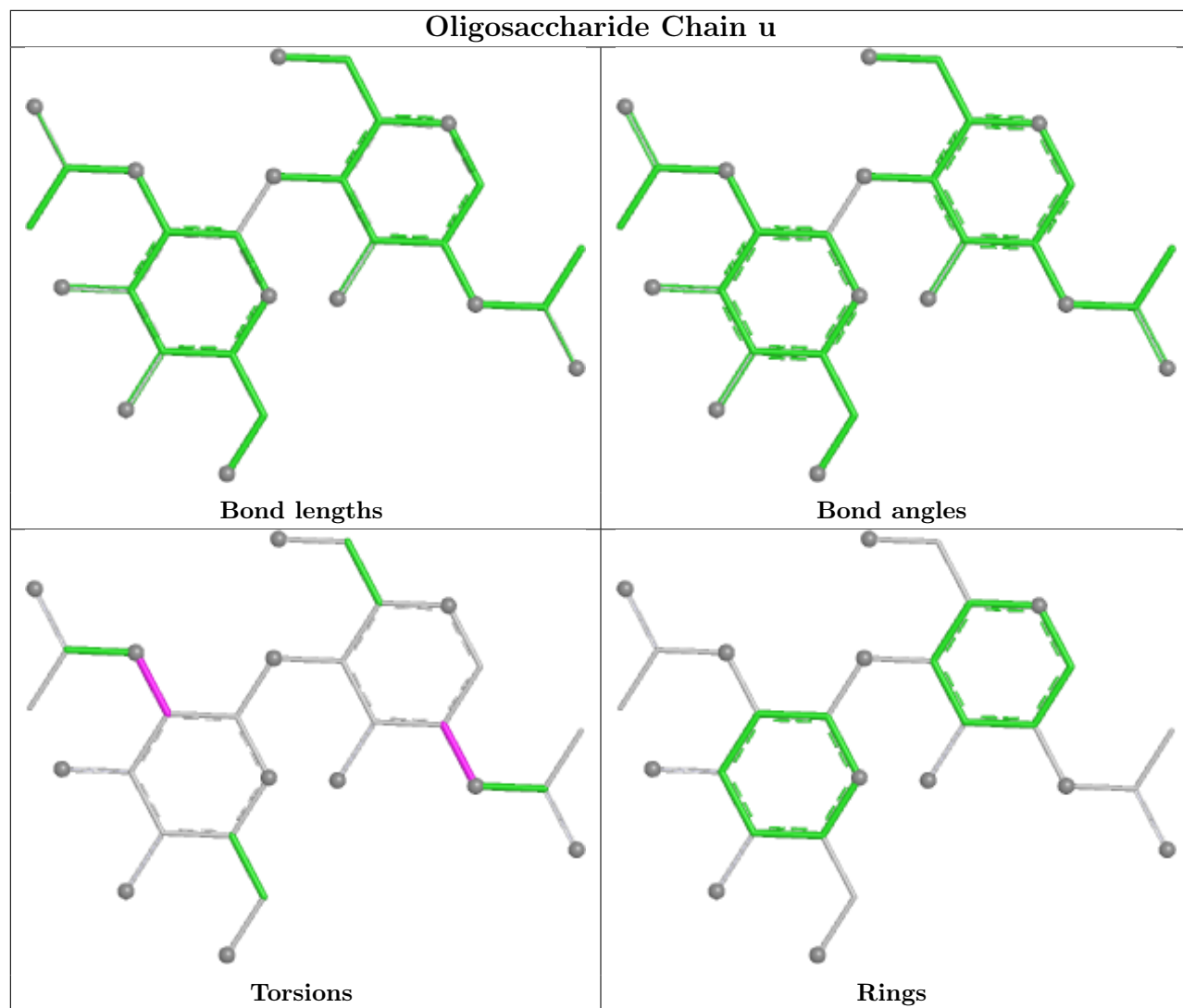


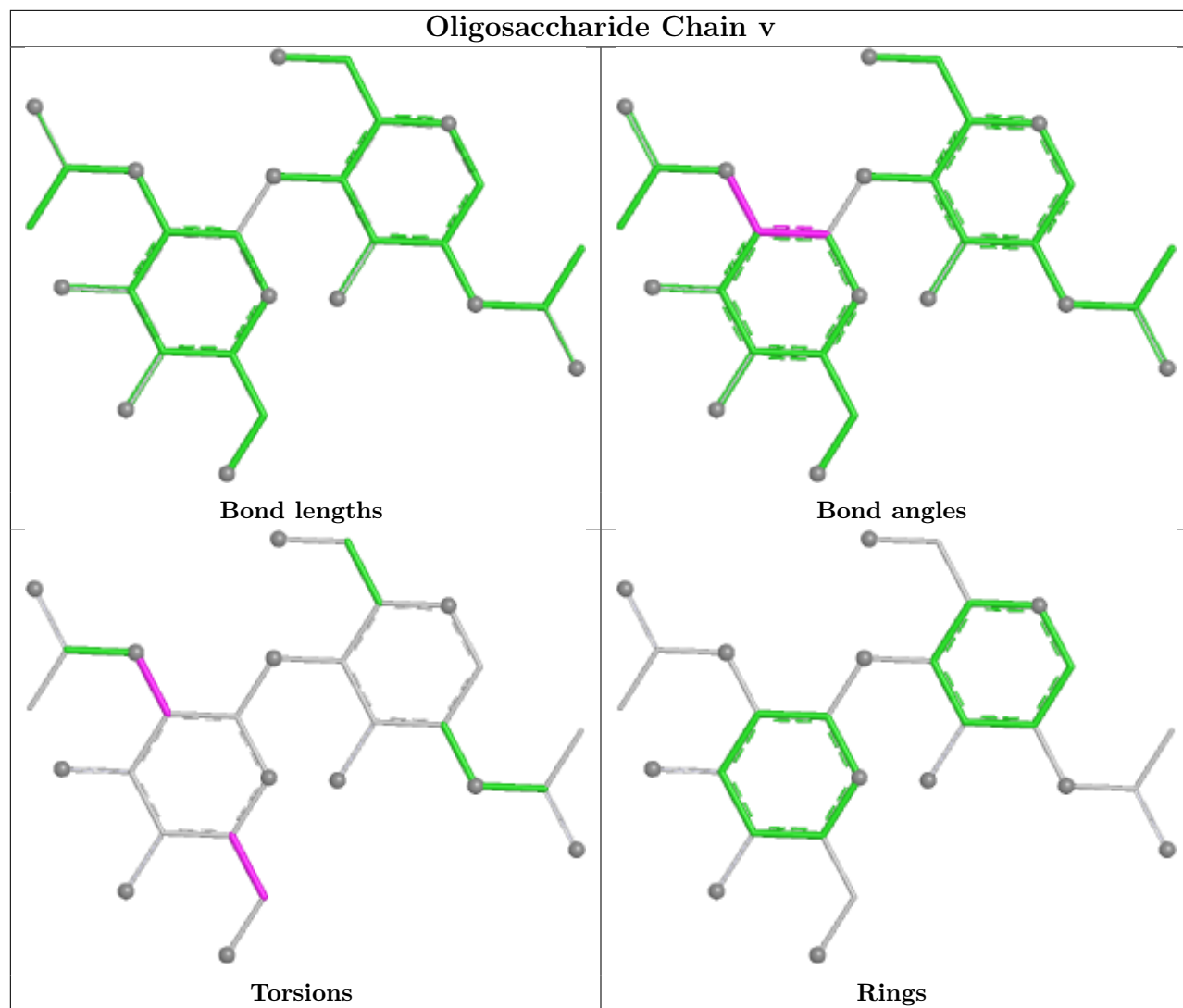


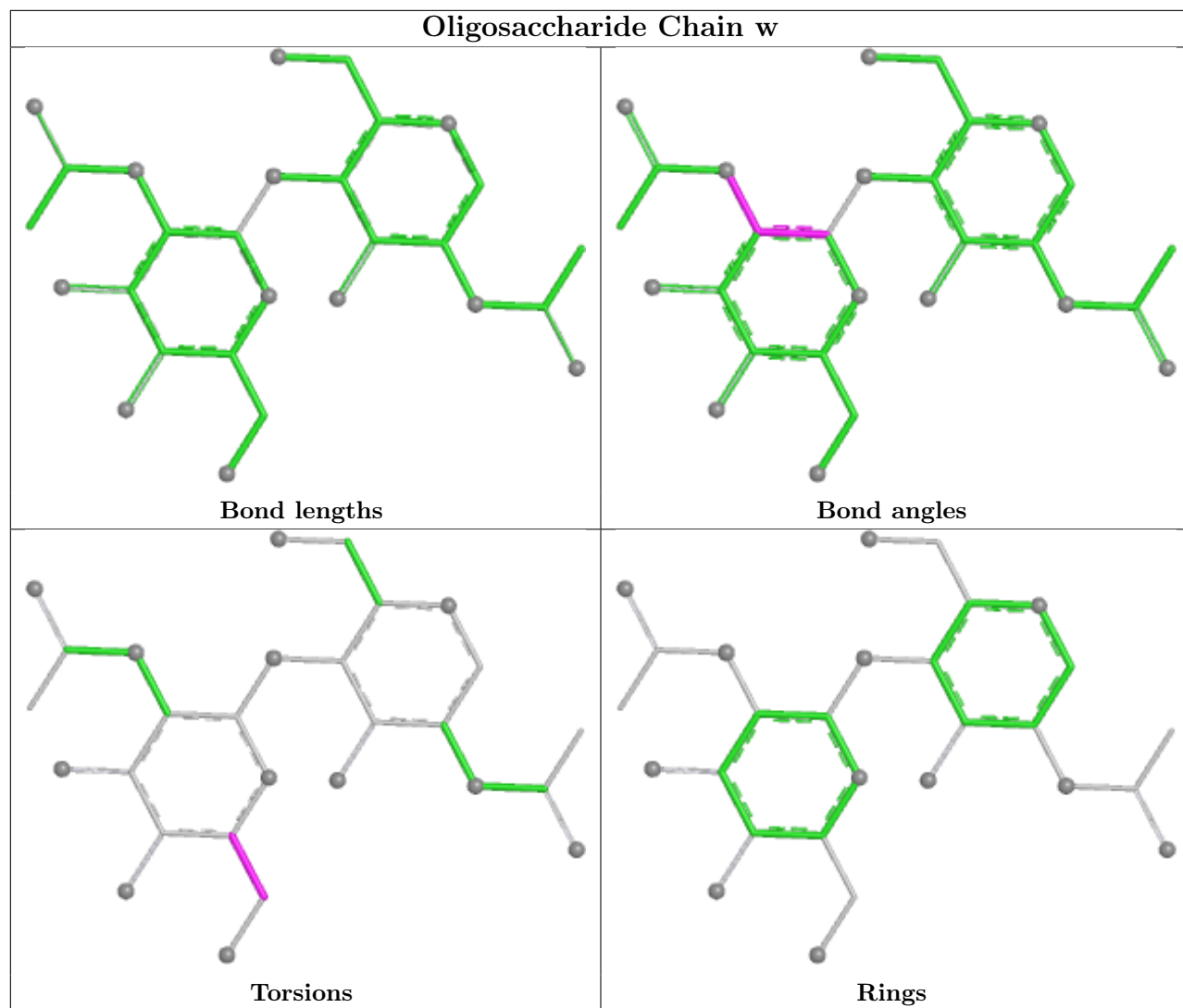


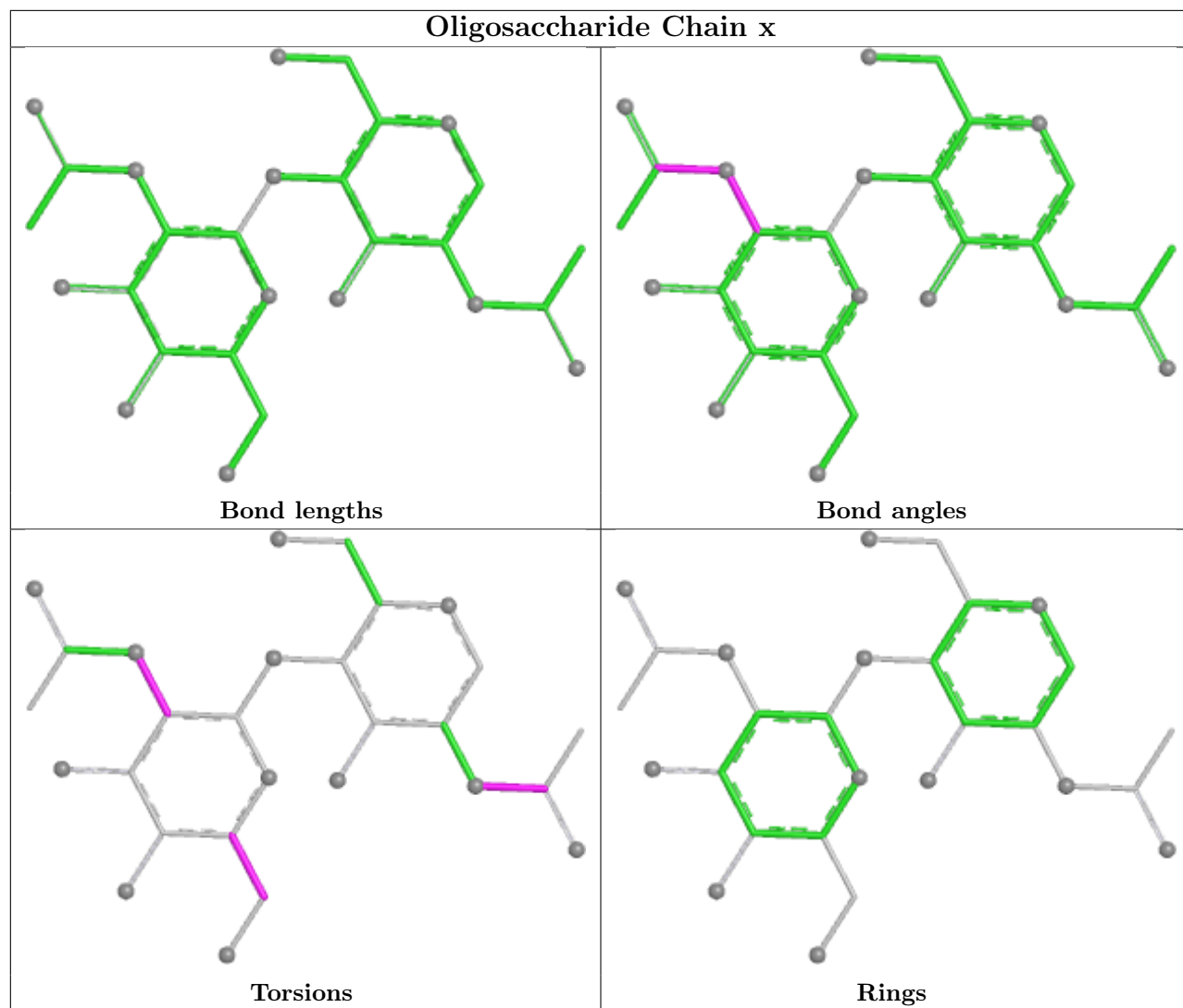


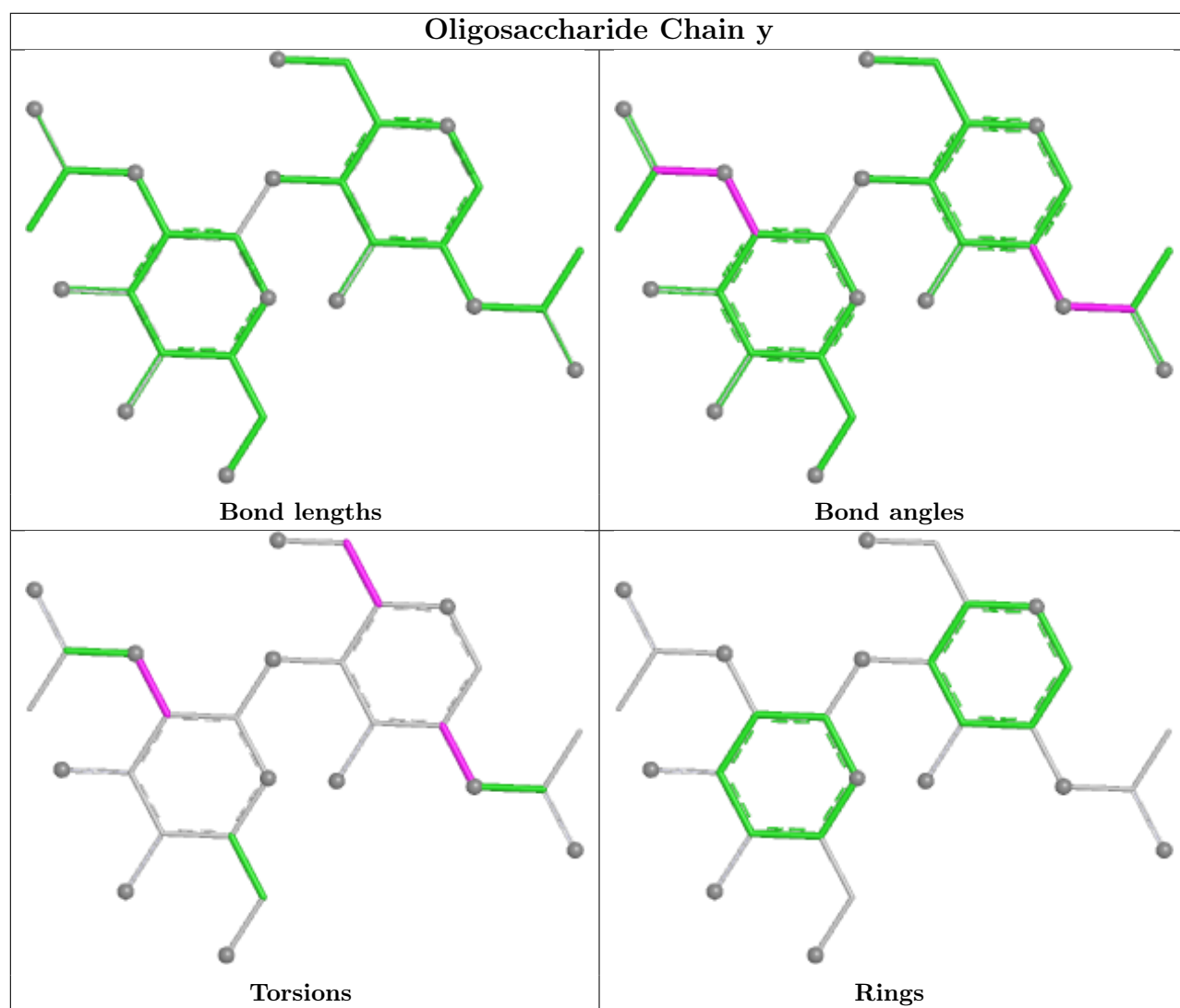


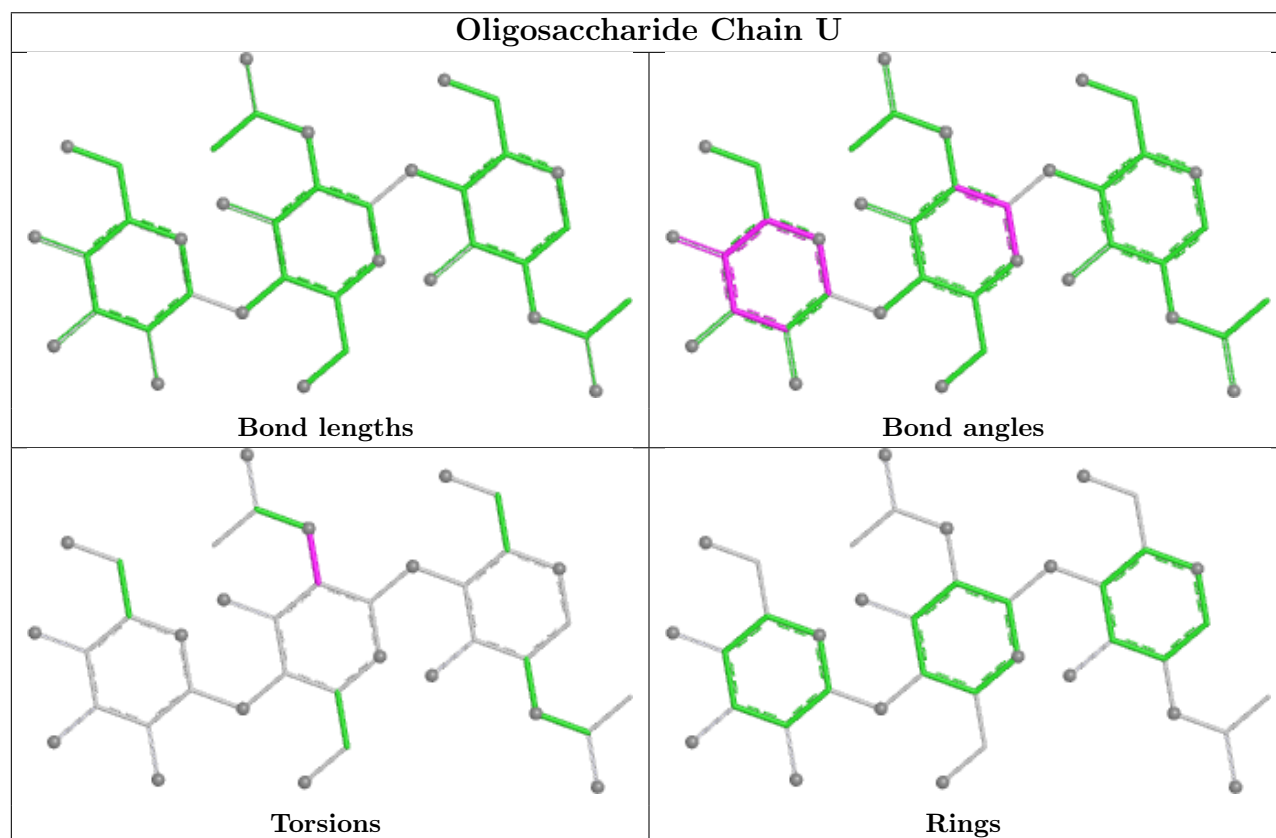
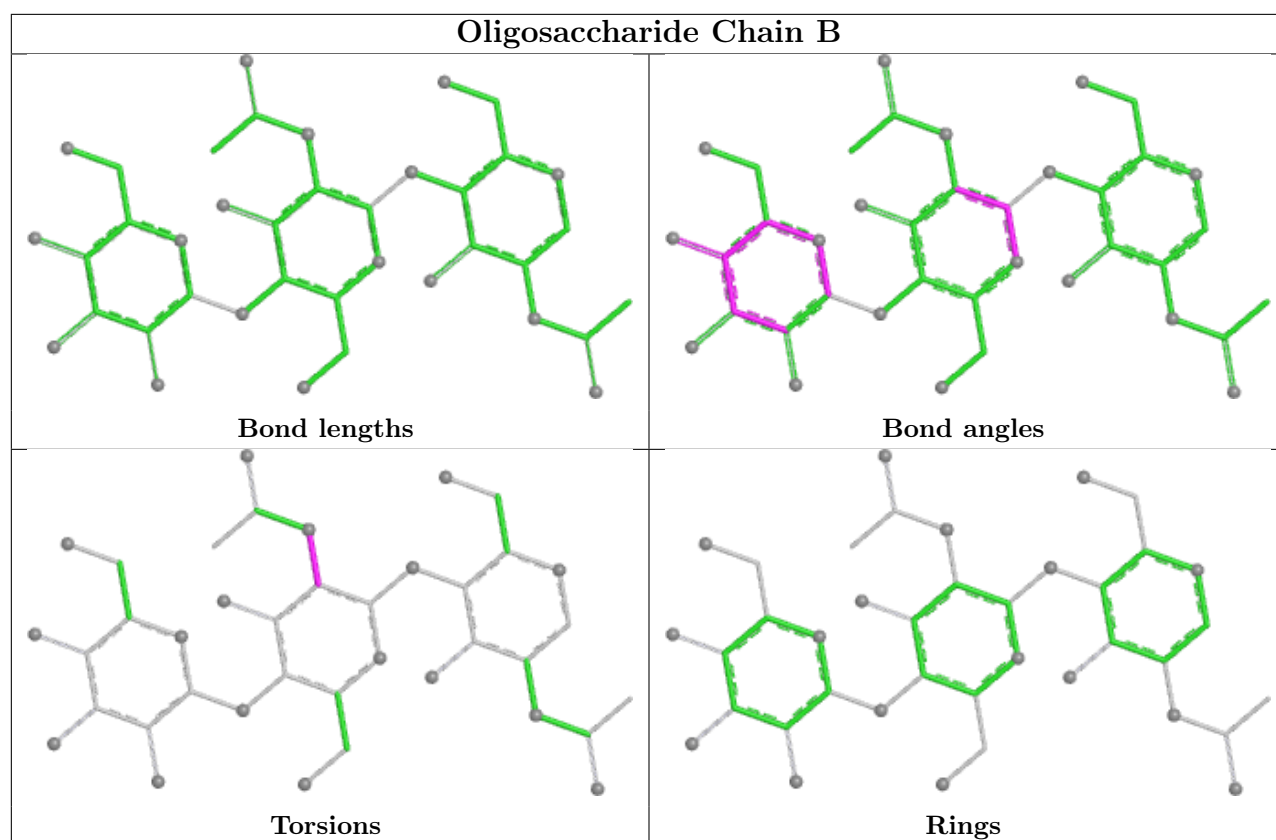


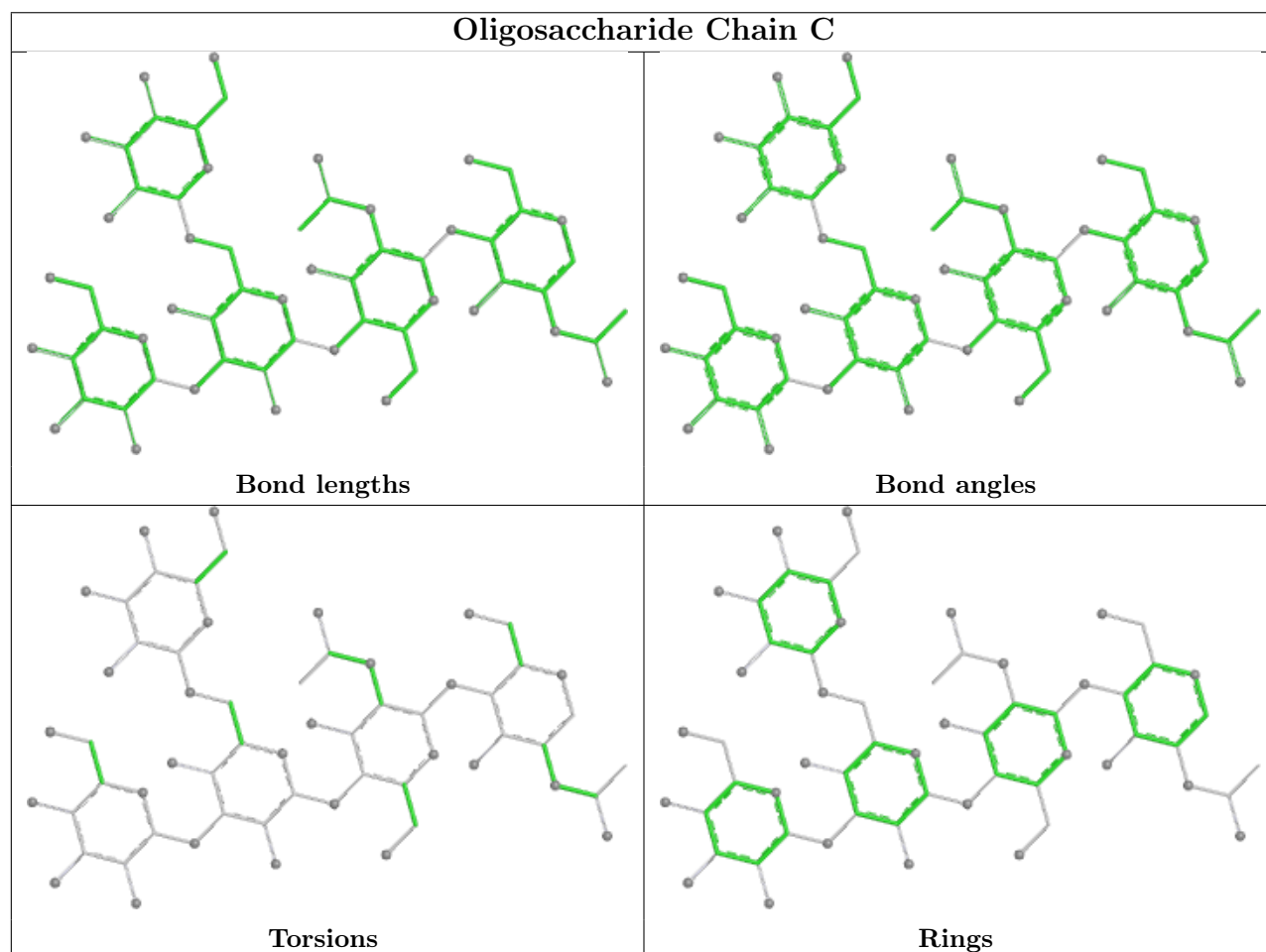
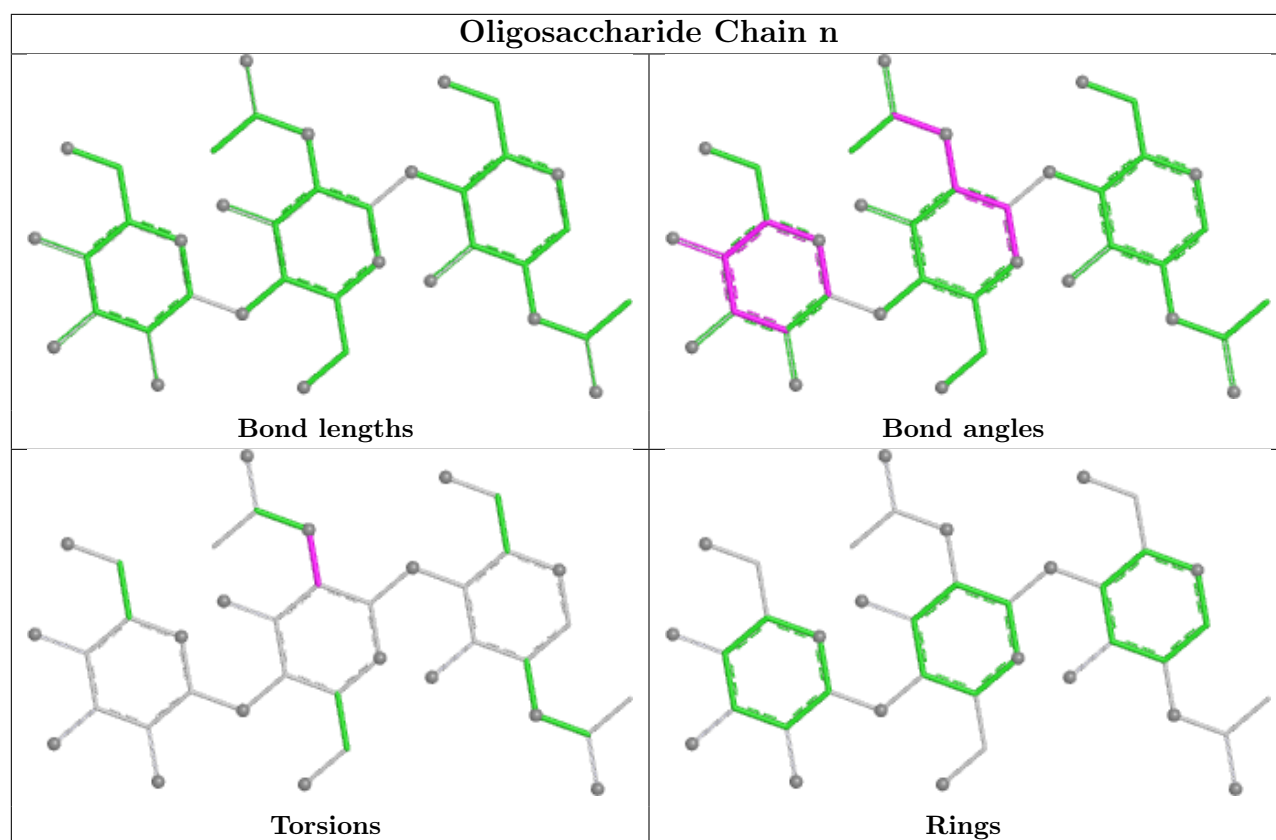


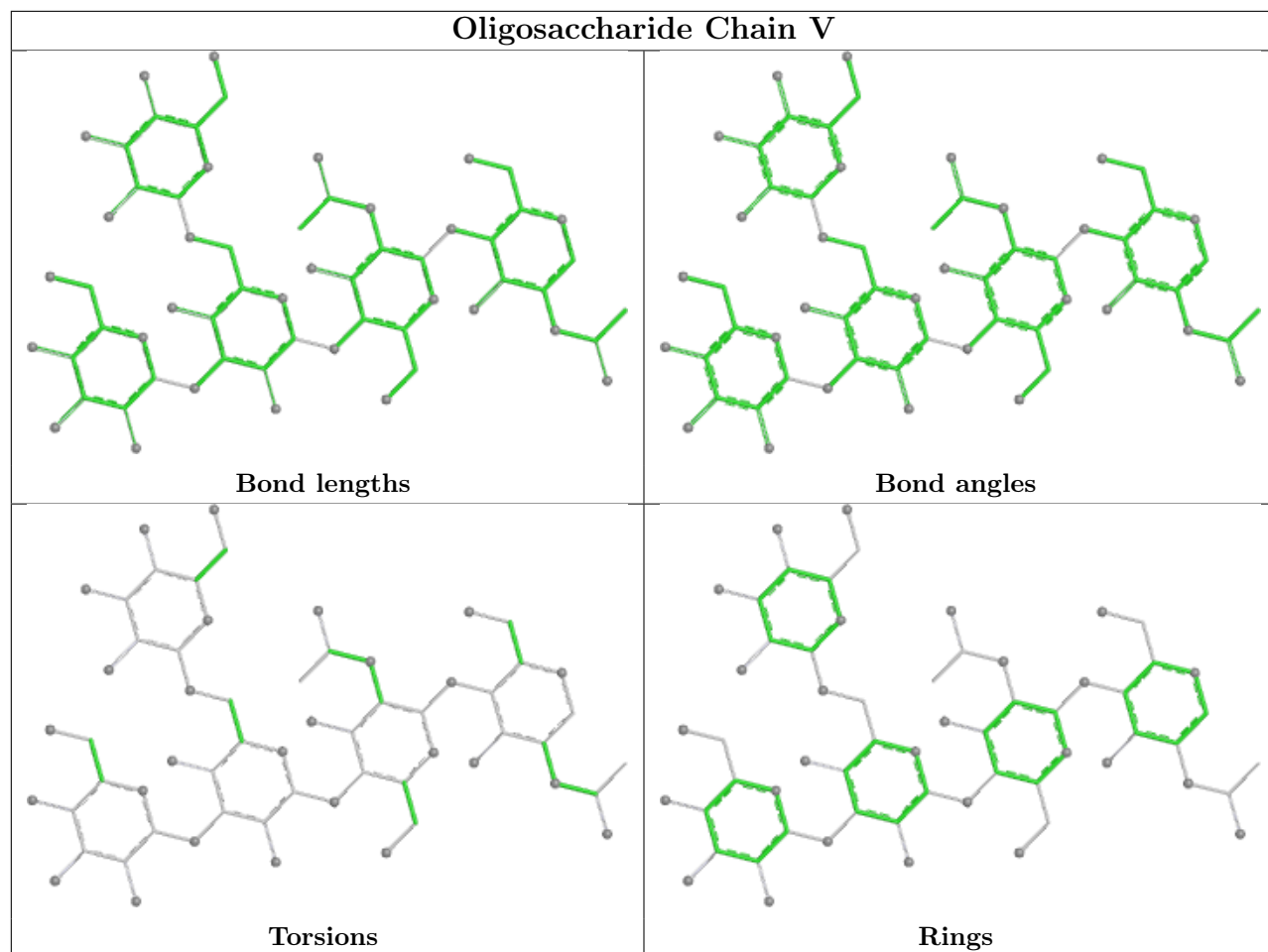


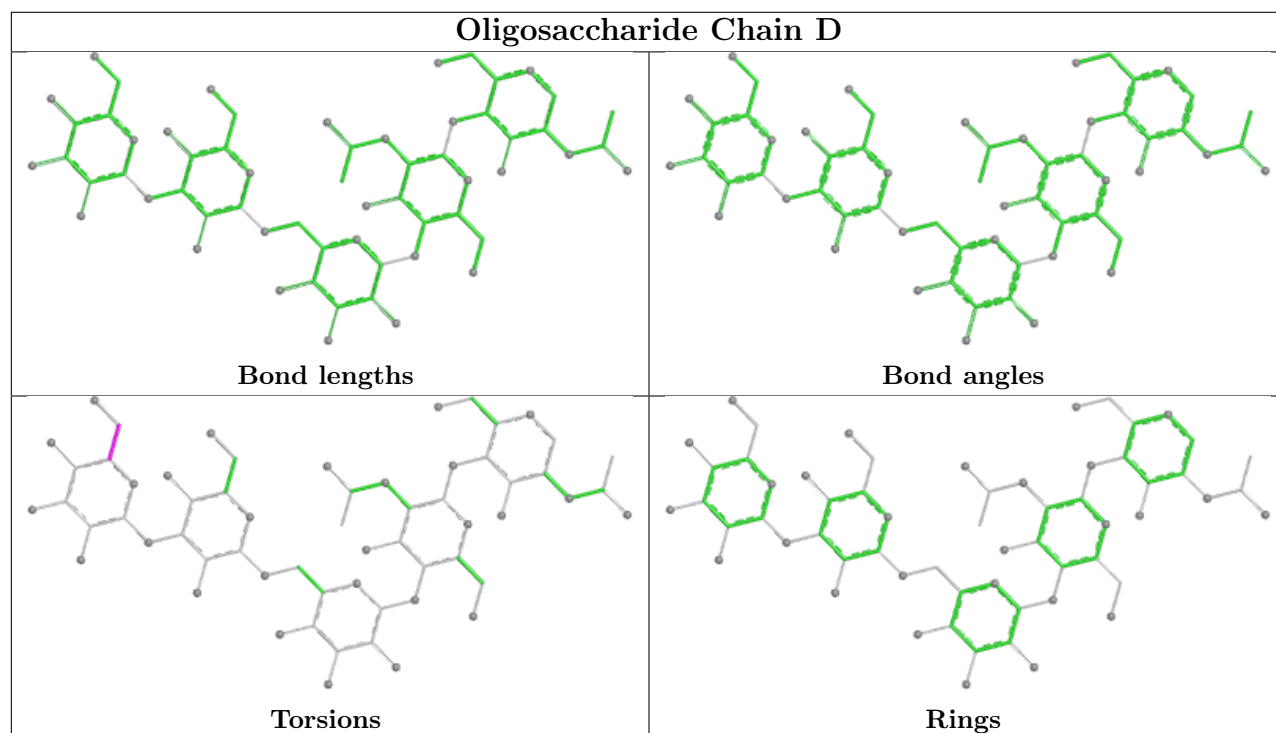
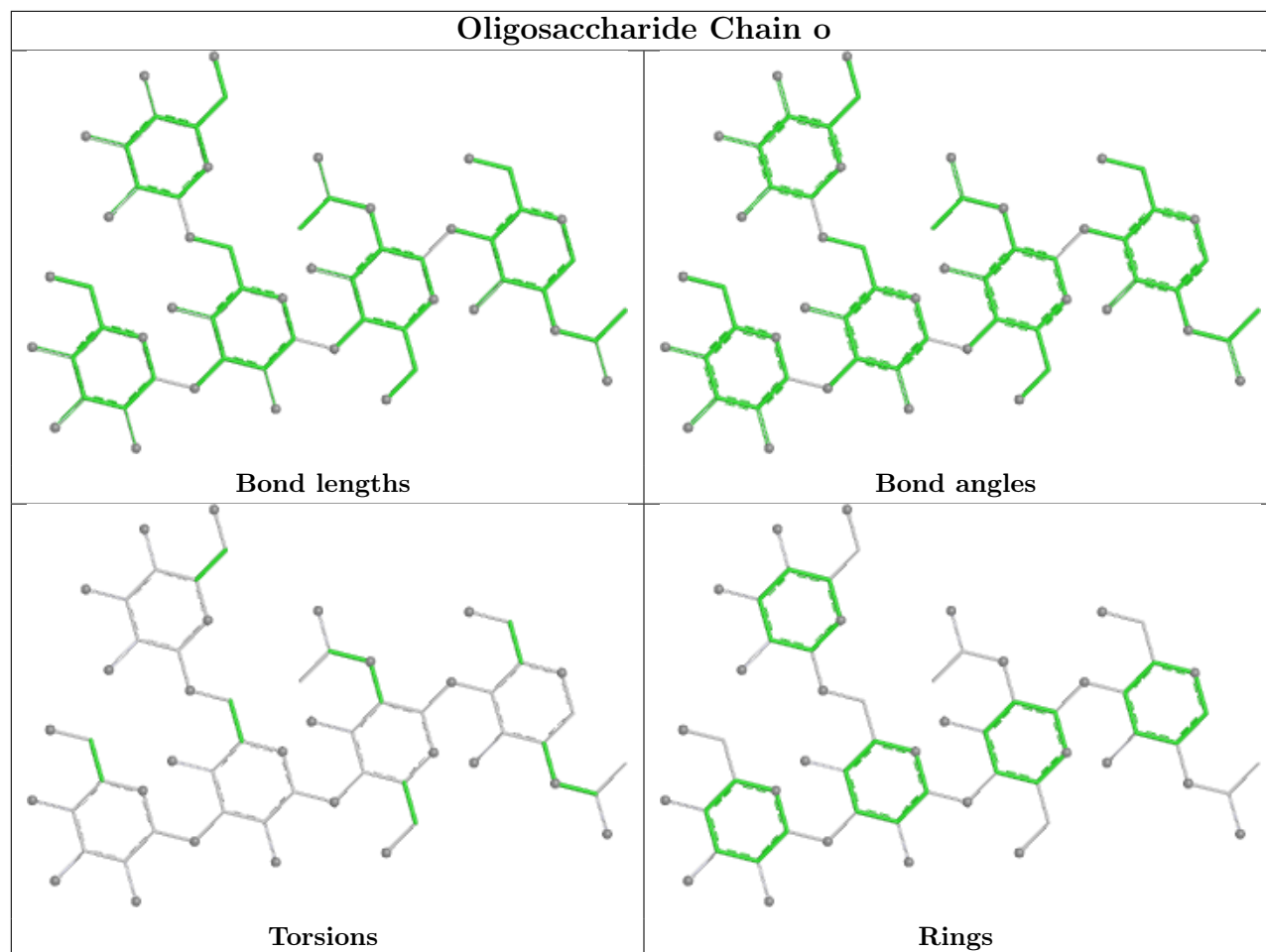


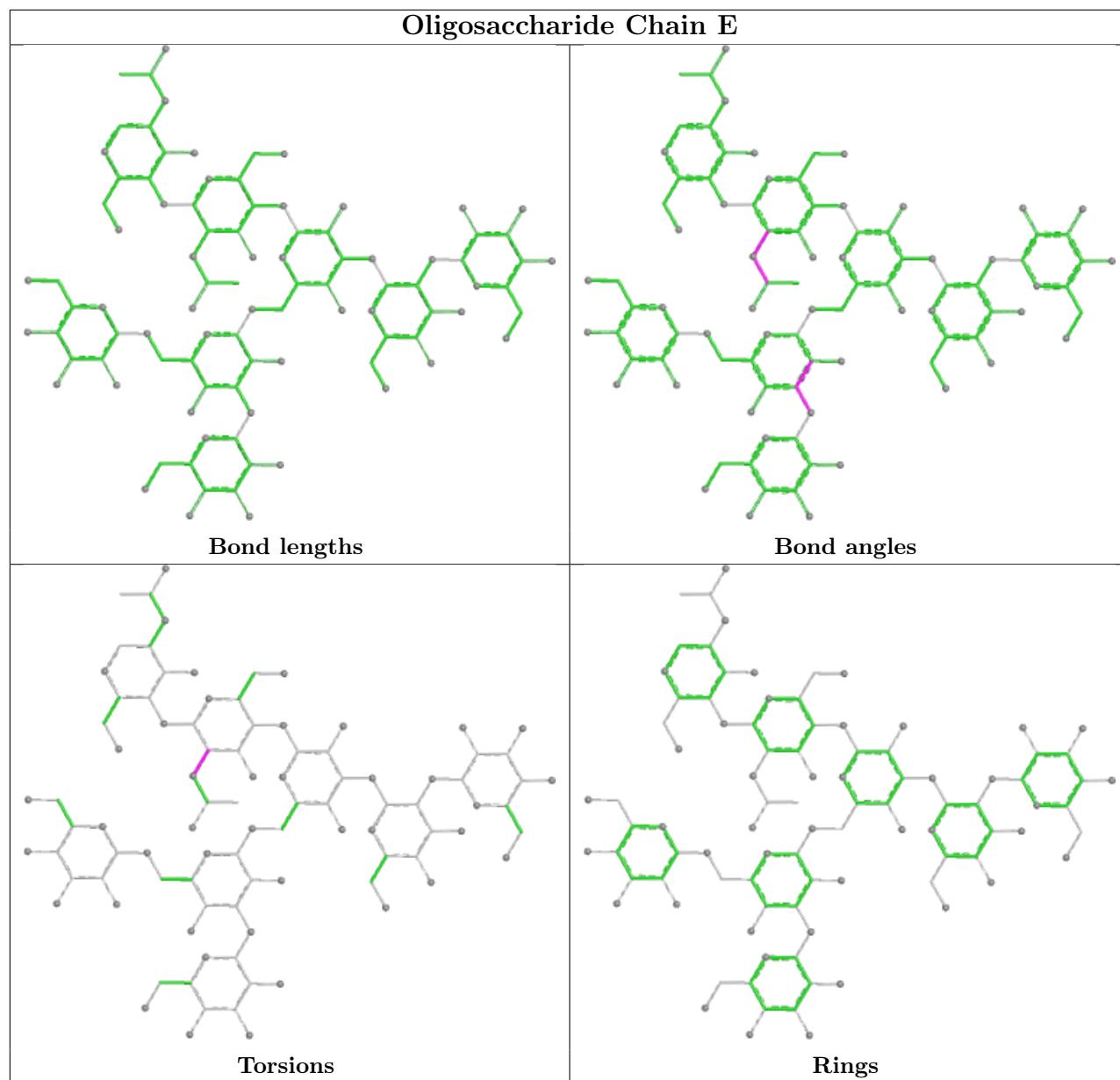


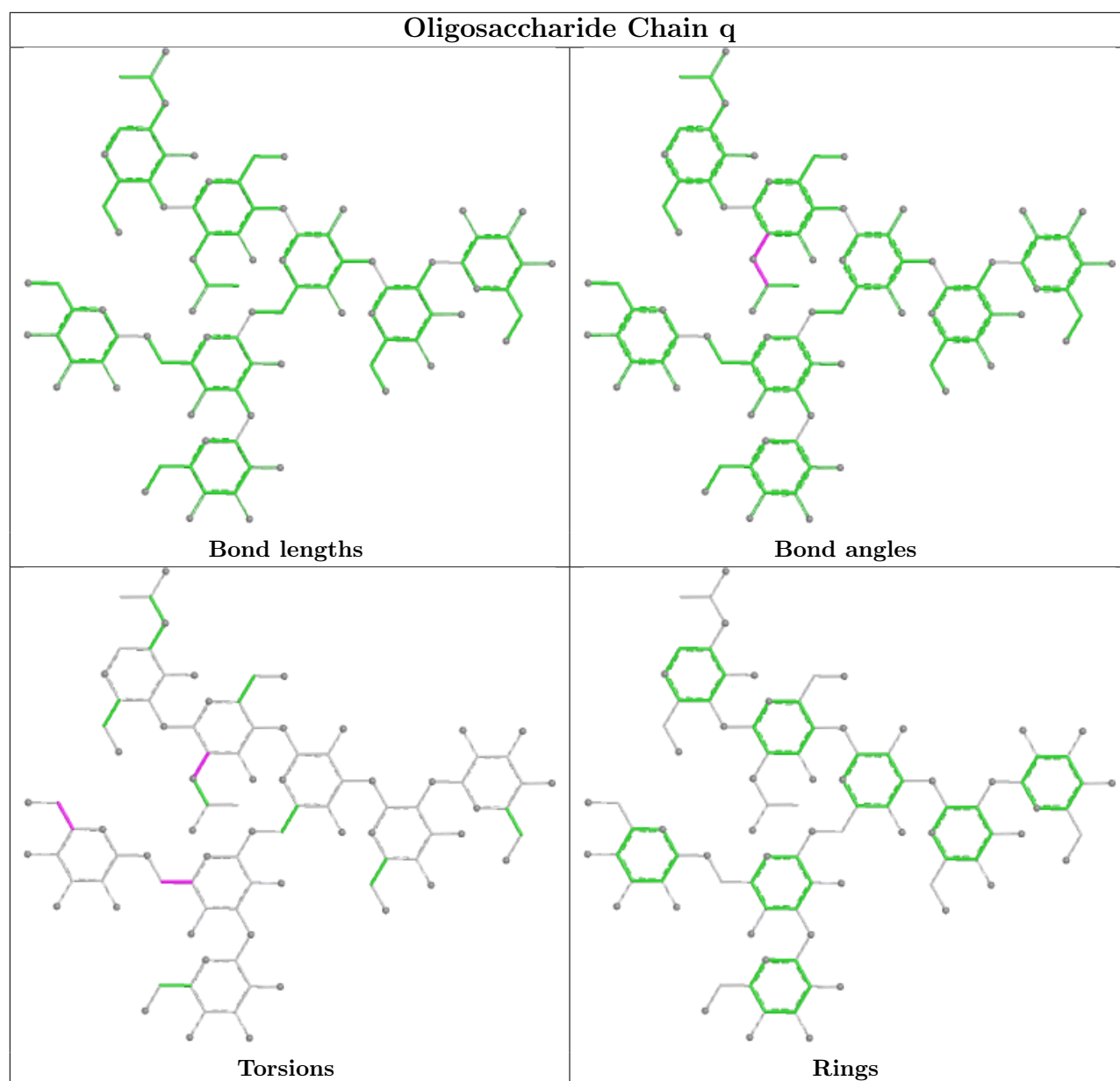


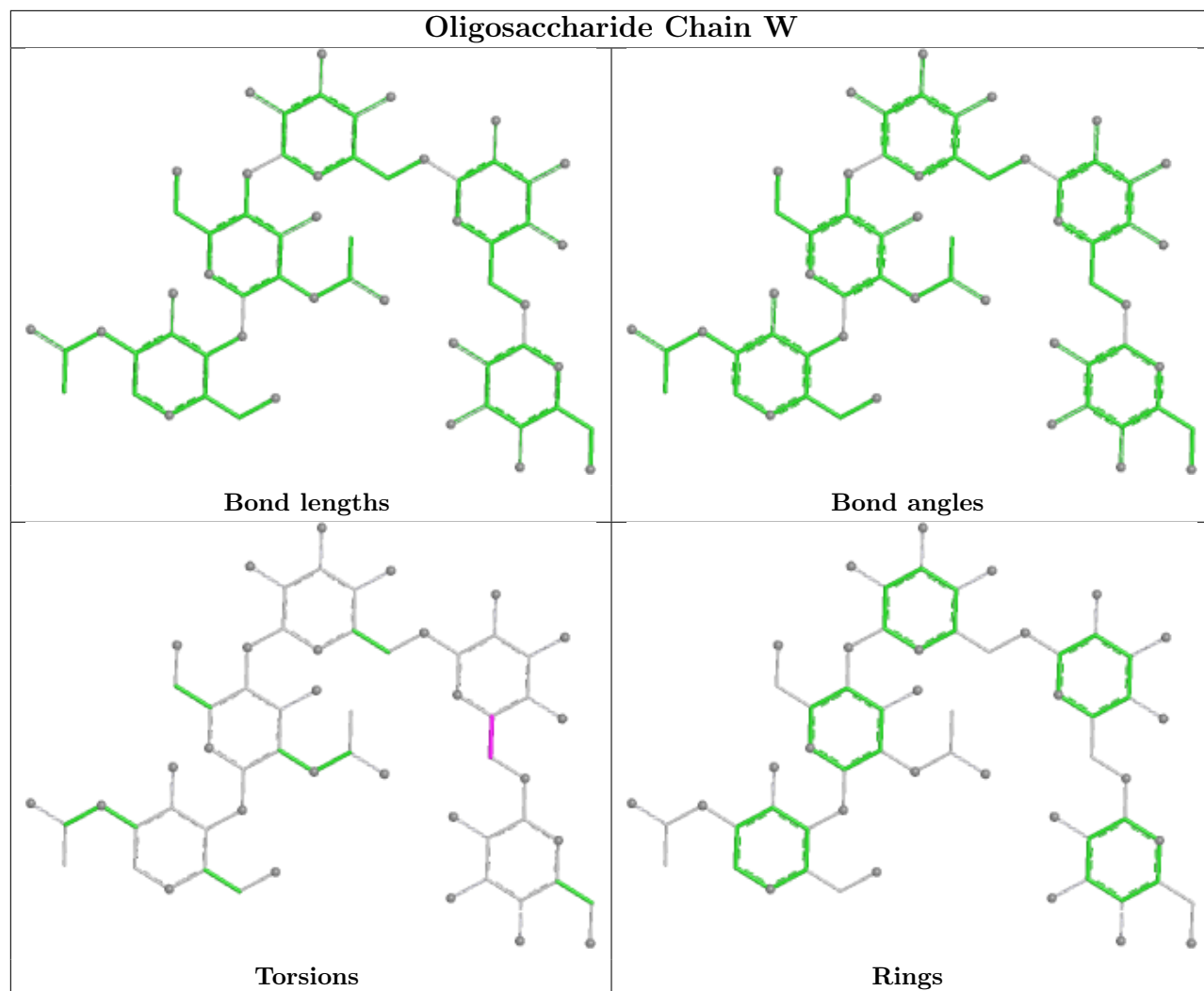


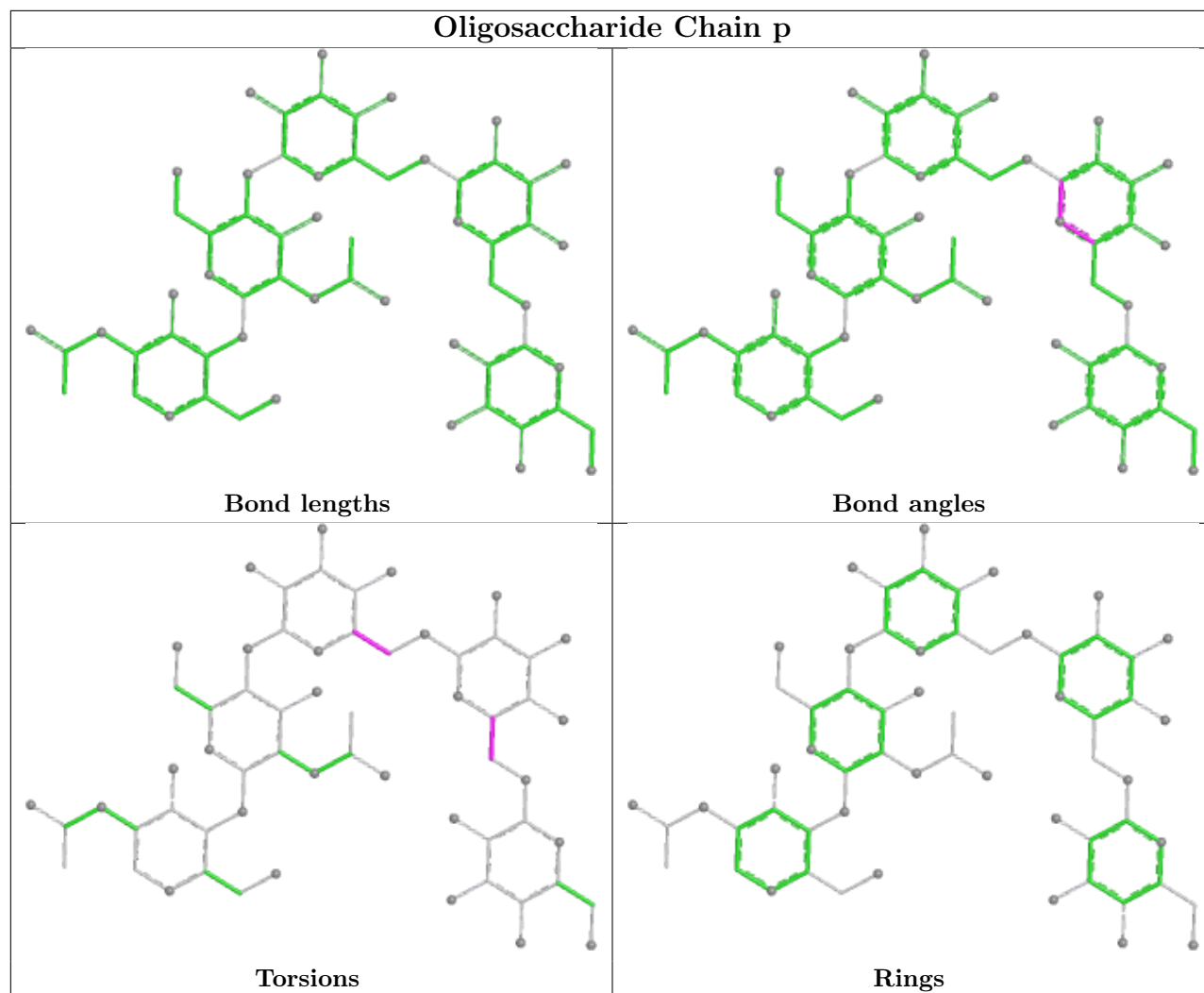


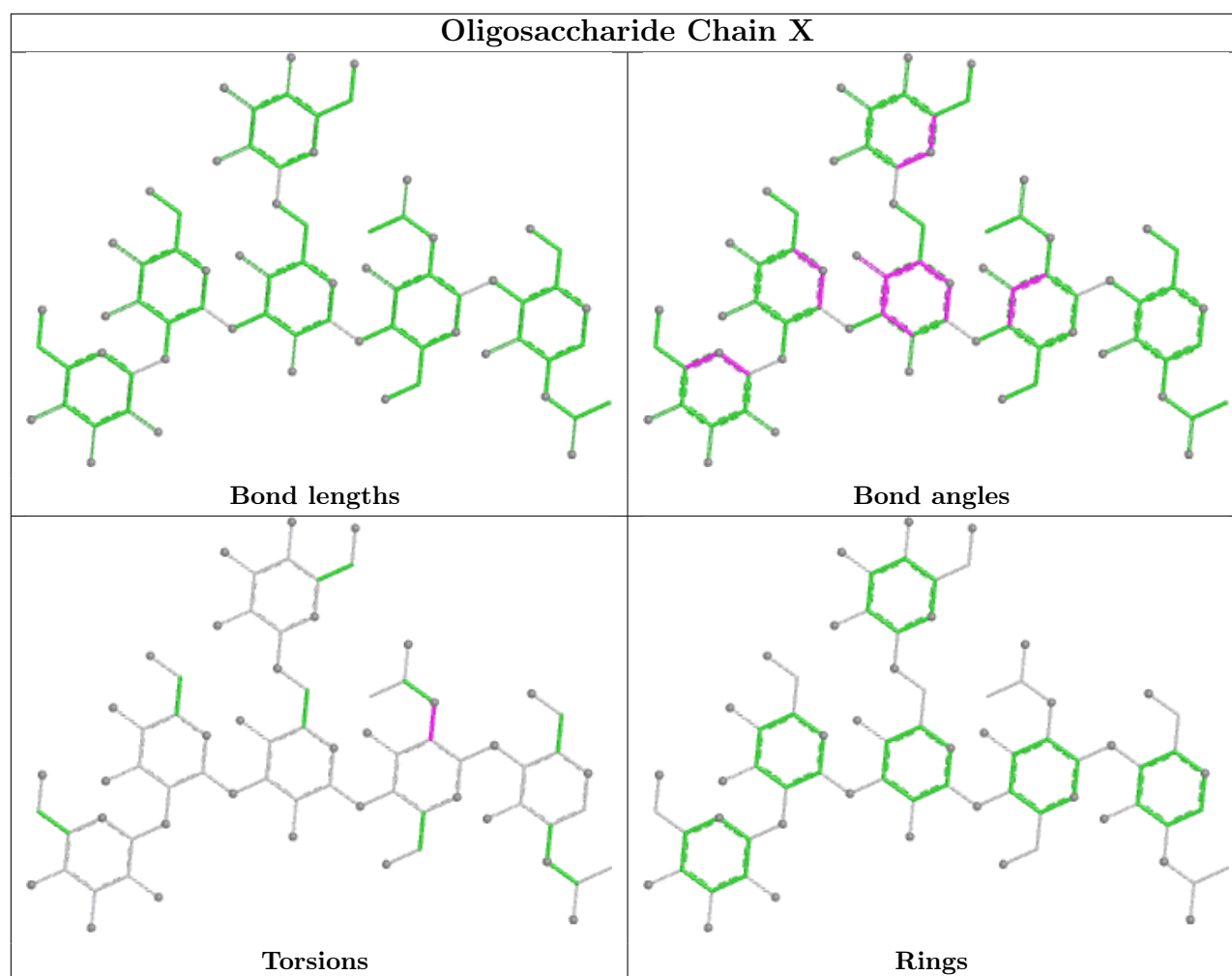












## 5.6 Ligand geometry [i](#)

30 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	$\# Z  > 2$
12	NAG	g	701	4	14,14,15	0.70	0	17,19,21	0.94	0
12	NAG	b	606	3	14,14,15	0.41	0	17,19,21	0.67	0
12	NAG	c	703	4	14,14,15	0.71	0	17,19,21	0.84	0
12	NAG	e	702	4	14,14,15	0.69	0	17,19,21	0.82	0
12	NAG	f	602	3	14,14,15	0.68	0	17,19,21	0.85	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
12	NAG	b	602	3	14,14,15	0.69	0	17,19,21	0.86	0
12	NAG	d	604	3	14,14,15	0.70	0	17,19,21	0.96	1 (5%)
12	NAG	b	604	3	14,14,15	0.37	0	17,19,21	0.47	0
12	NAG	f	607	3	14,14,15	0.39	0	17,19,21	0.92	2 (11%)
12	NAG	c	702	4	14,14,15	0.70	0	17,19,21	0.84	0
12	NAG	f	605	3	14,14,15	0.38	0	17,19,21	0.65	0
12	NAG	e	701	4	14,14,15	0.69	0	17,19,21	0.94	0
12	NAG	f	603	3	14,14,15	0.39	0	17,19,21	0.83	1 (5%)
12	NAG	g	703	4	14,14,15	0.70	0	17,19,21	0.84	0
12	NAG	c	701	4	14,14,15	0.71	0	17,19,21	0.93	0
12	NAG	e	703	4	14,14,15	0.69	0	17,19,21	0.83	0
12	NAG	b	603	3	14,14,15	0.40	0	17,19,21	0.85	1 (5%)
12	NAG	d	602	3	14,14,15	0.69	0	17,19,21	0.86	0
12	NAG	g	702	4	14,14,15	0.69	0	17,19,21	0.83	0
12	NAG	f	604	3	14,14,15	0.38	0	17,19,21	0.47	0
12	NAG	d	601	3	14,14,15	0.68	0	17,19,21	0.84	0
12	NAG	d	607	3	14,14,15	0.38	0	17,19,21	1.03	2 (11%)
12	NAG	b	605	3	14,14,15	0.39	0	17,19,21	0.66	0
12	NAG	d	605	3	14,14,15	0.71	0	17,19,21	0.84	0
12	NAG	b	607	3	14,14,15	0.38	0	17,19,21	0.48	0
12	NAG	f	601	3	14,14,15	0.68	0	17,19,21	0.83	0
12	NAG	d	606	3	14,14,15	0.39	0	17,19,21	0.53	0
12	NAG	b	601	3	14,14,15	0.68	0	17,19,21	0.84	0
12	NAG	f	606	3	14,14,15	0.40	0	17,19,21	1.00	2 (11%)
12	NAG	d	603	3	14,14,15	0.37	0	17,19,21	0.72	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	NAG	g	701	4	-	0/6/23/26	0/1/1/1
12	NAG	b	606	3	-	0/6/23/26	0/1/1/1
12	NAG	c	703	4	-	0/6/23/26	0/1/1/1
12	NAG	e	702	4	-	0/6/23/26	0/1/1/1
12	NAG	f	602	3	-	0/6/23/26	0/1/1/1
12	NAG	b	602	3	-	0/6/23/26	0/1/1/1
12	NAG	d	604	3	-	0/6/23/26	0/1/1/1
12	NAG	b	604	3	-	0/6/23/26	0/1/1/1
12	NAG	f	607	3	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	NAG	c	702	4	-	0/6/23/26	0/1/1/1
12	NAG	f	605	3	-	0/6/23/26	0/1/1/1
12	NAG	e	701	4	-	0/6/23/26	0/1/1/1
12	NAG	f	603	3	-	3/6/23/26	0/1/1/1
12	NAG	g	703	4	-	0/6/23/26	0/1/1/1
12	NAG	c	701	4	-	0/6/23/26	0/1/1/1
12	NAG	e	703	4	-	0/6/23/26	0/1/1/1
12	NAG	b	603	3	-	3/6/23/26	0/1/1/1
12	NAG	d	602	3	-	0/6/23/26	0/1/1/1
12	NAG	g	702	4	-	0/6/23/26	0/1/1/1
12	NAG	f	604	3	-	0/6/23/26	0/1/1/1
12	NAG	d	601	3	-	0/6/23/26	0/1/1/1
12	NAG	d	607	3	-	4/6/23/26	0/1/1/1
12	NAG	b	605	3	-	0/6/23/26	0/1/1/1
12	NAG	d	605	3	-	0/6/23/26	0/1/1/1
12	NAG	b	607	3	-	1/6/23/26	0/1/1/1
12	NAG	f	601	3	-	0/6/23/26	0/1/1/1
12	NAG	d	606	3	-	0/6/23/26	0/1/1/1
12	NAG	b	601	3	-	0/6/23/26	0/1/1/1
12	NAG	f	606	3	-	4/6/23/26	0/1/1/1
12	NAG	d	603	3	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	d	607	NAG	C1-C2-N2	3.16	115.41	110.43
12	f	606	NAG	C1-C2-N2	3.08	115.29	110.43
12	f	607	NAG	C1-C2-N2	2.43	114.25	110.43
12	f	603	NAG	C1-C2-N2	2.42	114.25	110.43
12	b	603	NAG	C1-C2-N2	2.36	114.15	110.43
12	f	606	NAG	C2-N2-C7	2.18	125.82	122.90
12	d	607	NAG	C2-N2-C7	2.11	125.73	122.90
12	d	604	NAG	C1-O5-C5	2.06	114.95	112.19
12	f	607	NAG	C1-O5-C5	2.00	114.87	112.19

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	f	607	NAG	C1-C2-N2-C7
12	f	607	NAG	C8-C7-N2-C2
12	f	607	NAG	O7-C7-N2-C2
12	f	603	NAG	C8-C7-N2-C2
12	d	607	NAG	C8-C7-N2-C2
12	b	603	NAG	C8-C7-N2-C2
12	f	603	NAG	O7-C7-N2-C2
12	b	607	NAG	O5-C5-C6-O6
12	d	607	NAG	O7-C7-N2-C2
12	b	603	NAG	C3-C2-N2-C7
12	d	603	NAG	C3-C2-N2-C7
12	d	607	NAG	C3-C2-N2-C7
12	f	603	NAG	C3-C2-N2-C7
12	f	606	NAG	C3-C2-N2-C7
12	f	607	NAG	C3-C2-N2-C7
12	f	606	NAG	C8-C7-N2-C2
12	d	607	NAG	C1-C2-N2-C7
12	f	606	NAG	C1-C2-N2-C7
12	b	603	NAG	O7-C7-N2-C2
12	f	606	NAG	O7-C7-N2-C2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	f	607	NAG	1	0
12	b	603	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.

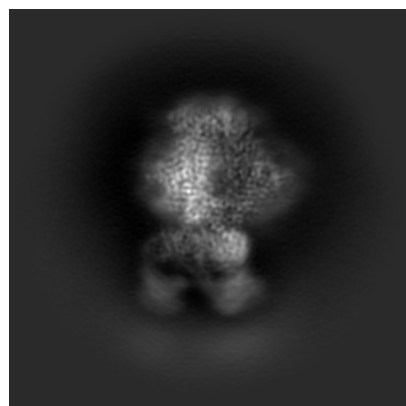
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-70666. These allow visual inspection of the internal detail of the map and identification of artifacts.

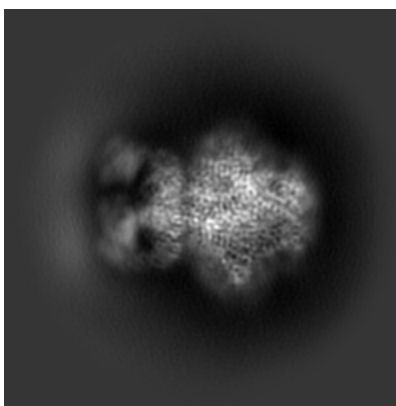
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

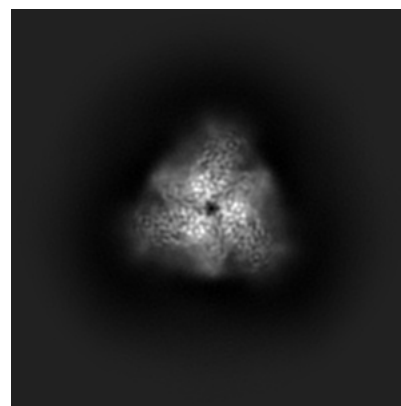
#### 6.1.1 Primary map



X

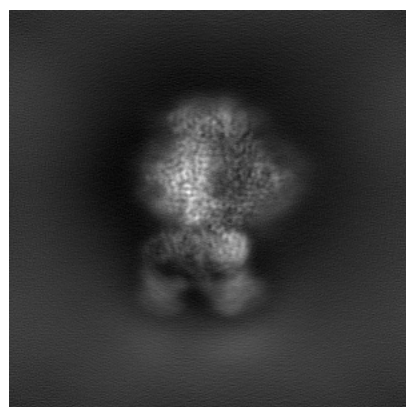


Y

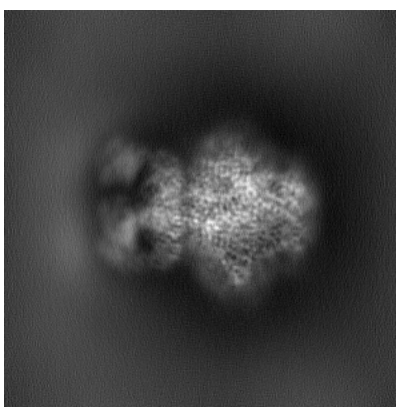


Z

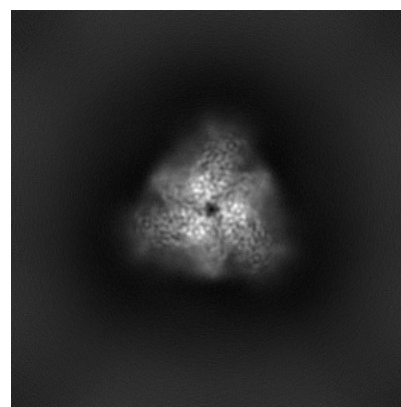
#### 6.1.2 Raw map



X



Y

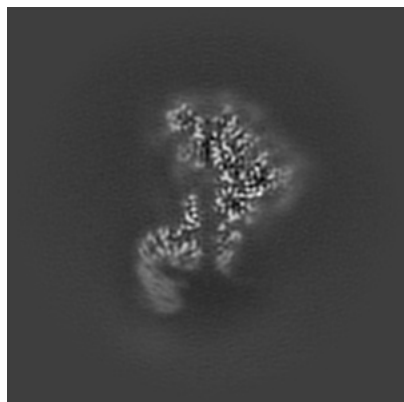


Z

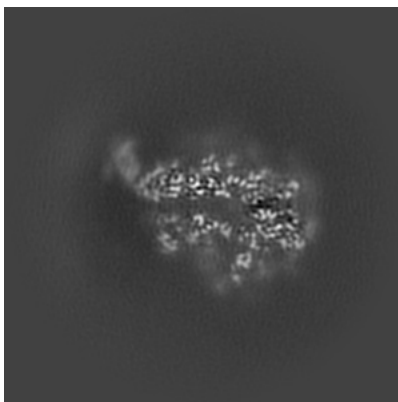
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

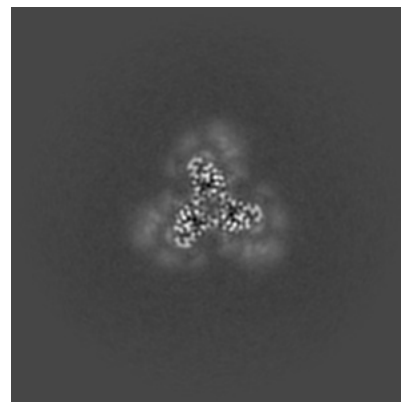
### 6.2.1 Primary map



X Index: 192

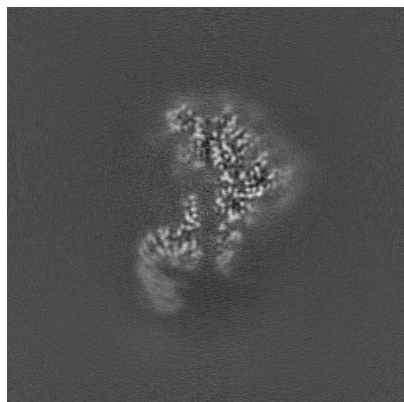


Y Index: 192

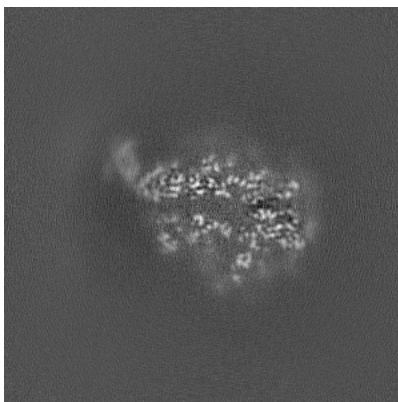


Z Index: 192

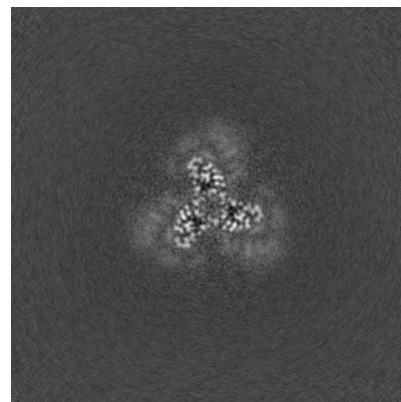
### 6.2.2 Raw map



X Index: 192



Y Index: 192

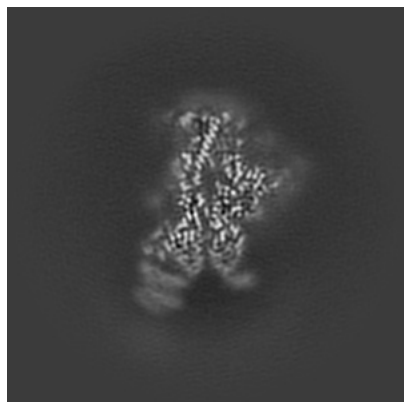


Z Index: 192

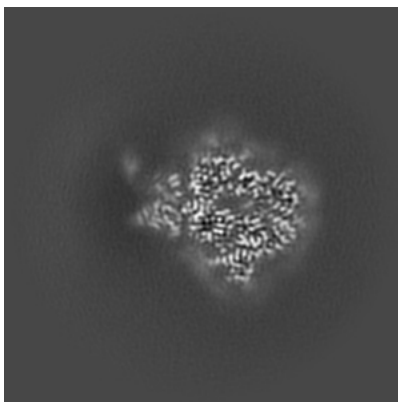
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

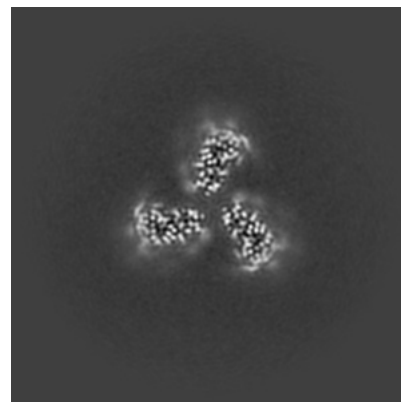
### 6.3.1 Primary map



X Index: 183

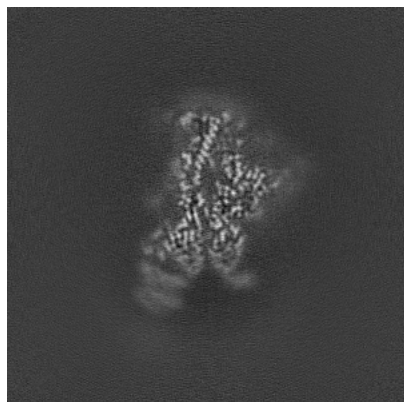


Y Index: 180

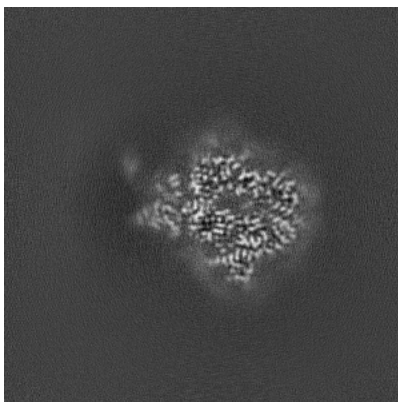


Z Index: 223

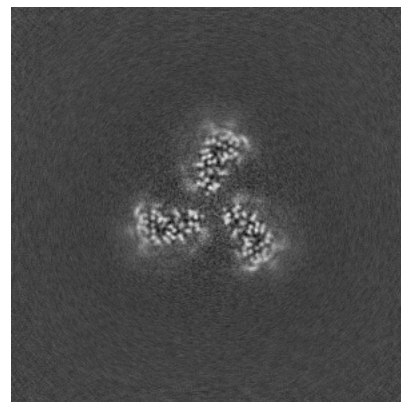
### 6.3.2 Raw map



X Index: 183



Y Index: 180

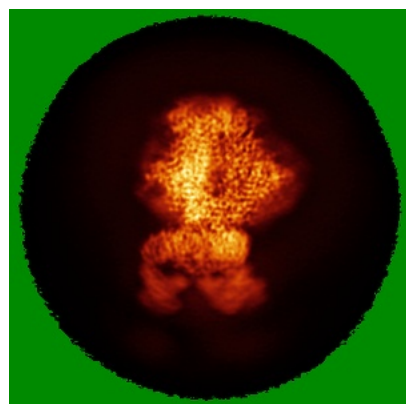


Z Index: 223

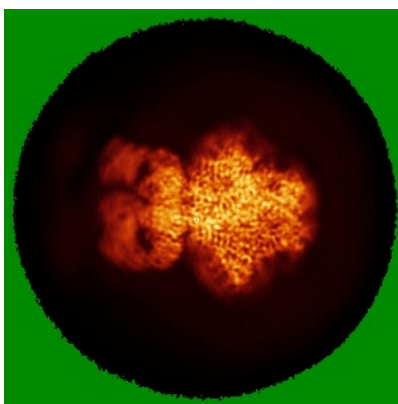
The images above show the largest variance slices of the map in three orthogonal directions.

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

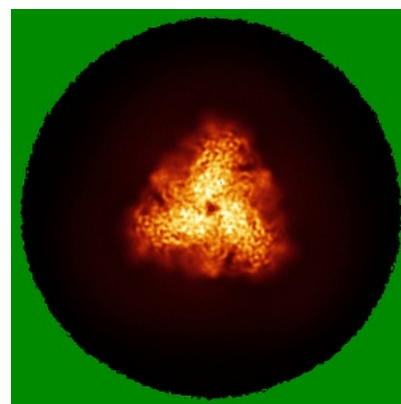
### 6.4.1 Primary map



X

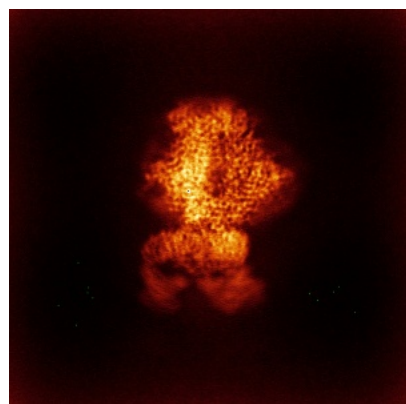


Y

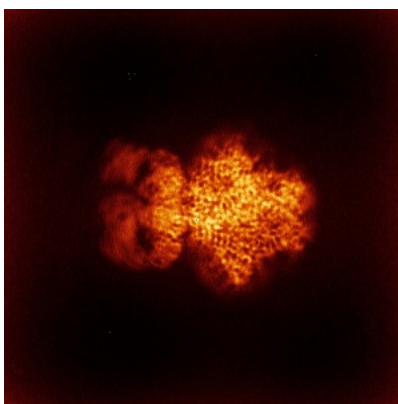


Z

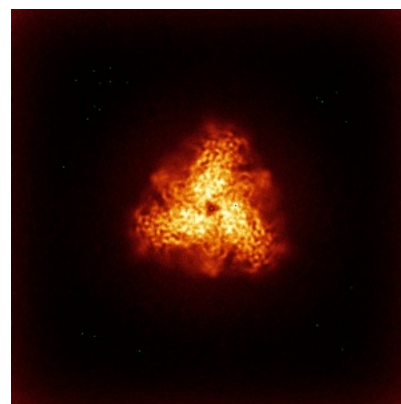
### 6.4.2 Raw map



X



Y

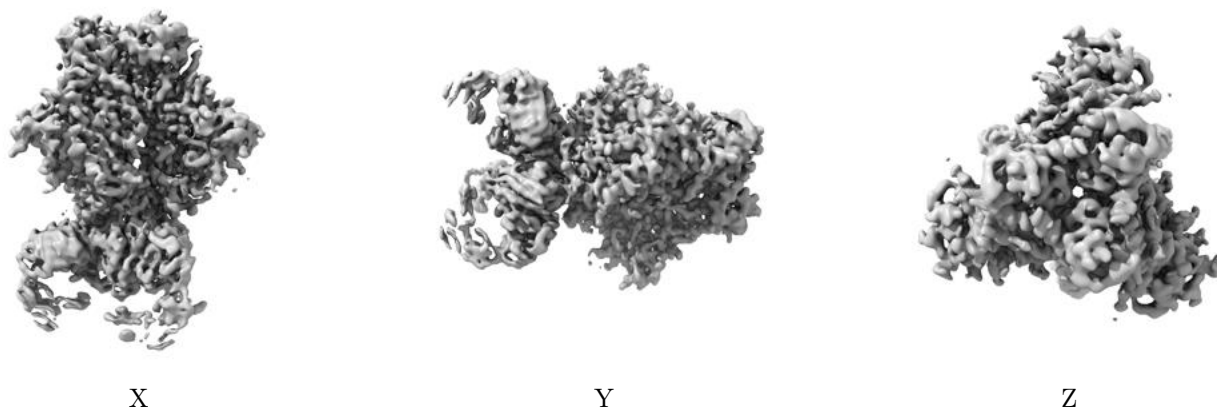


Z

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

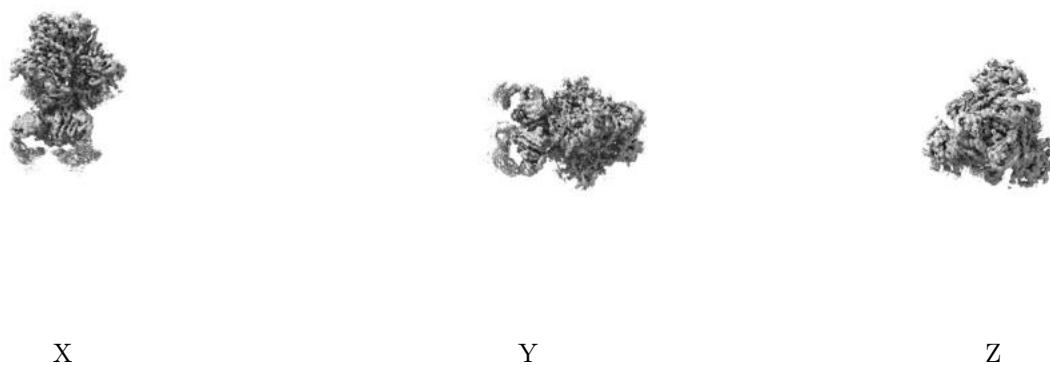
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.09. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.5.2 Raw map



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

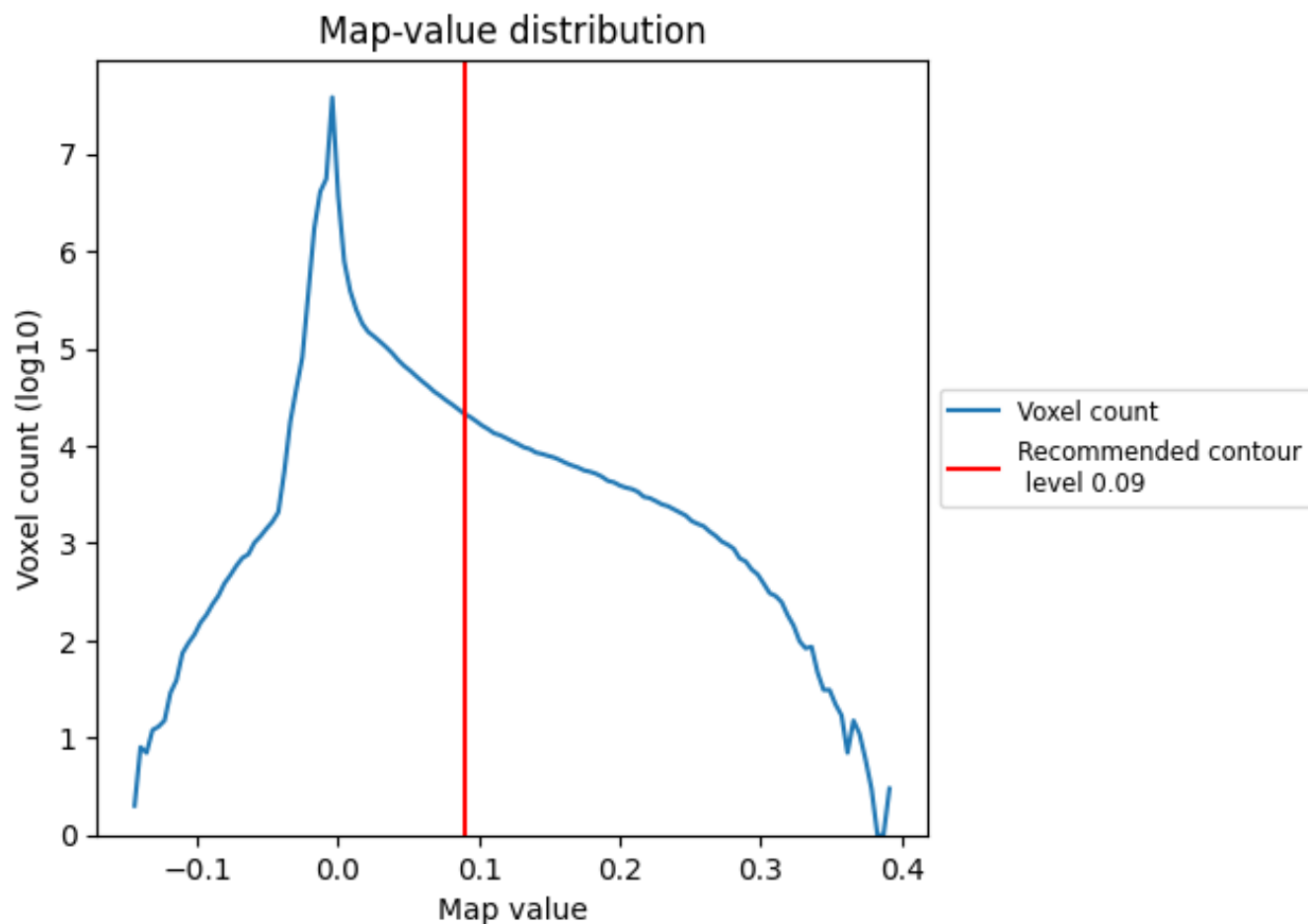
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

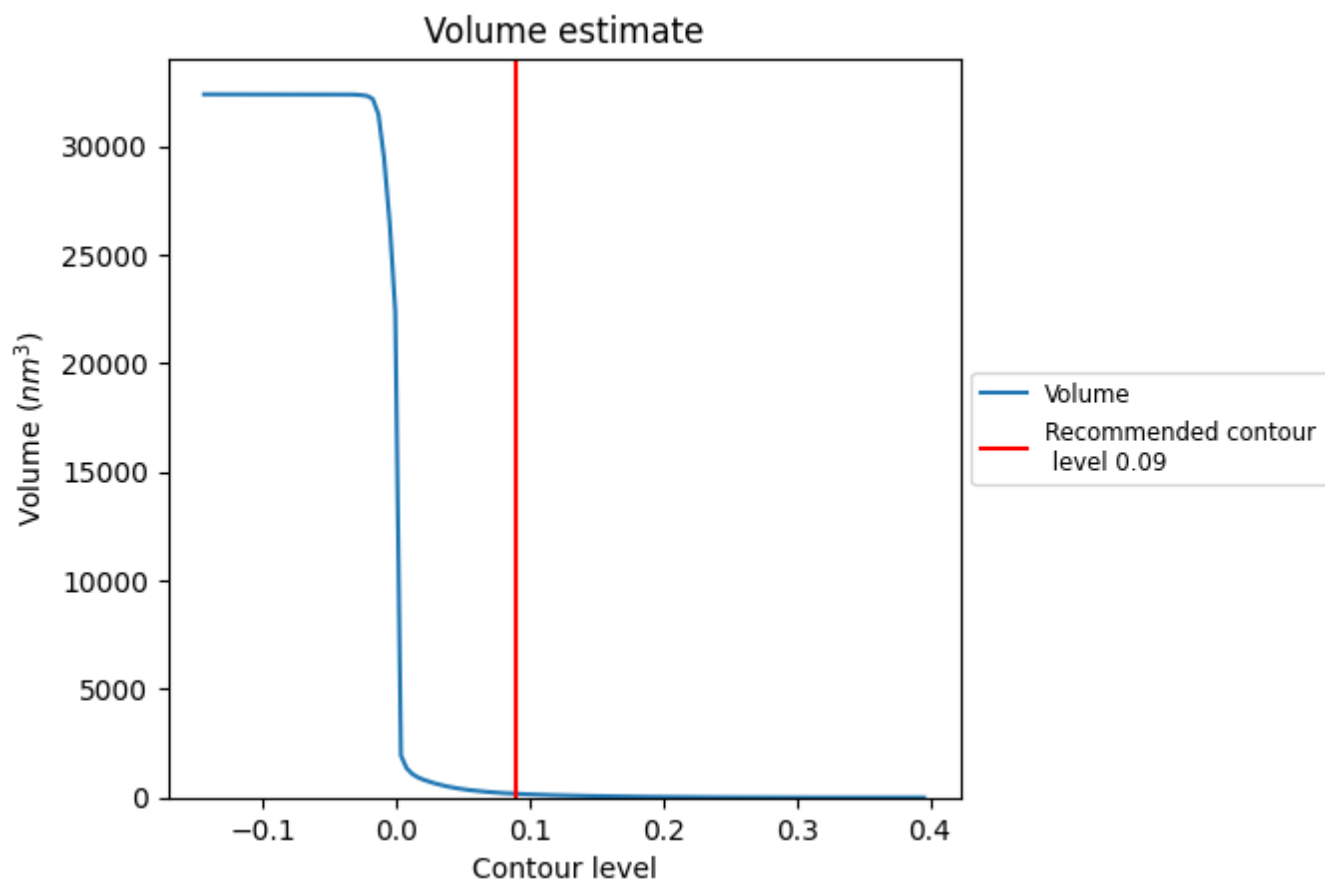
This section contains the results of statistical analysis of the map.

### 7.1 Map-value distribution [i](#)



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

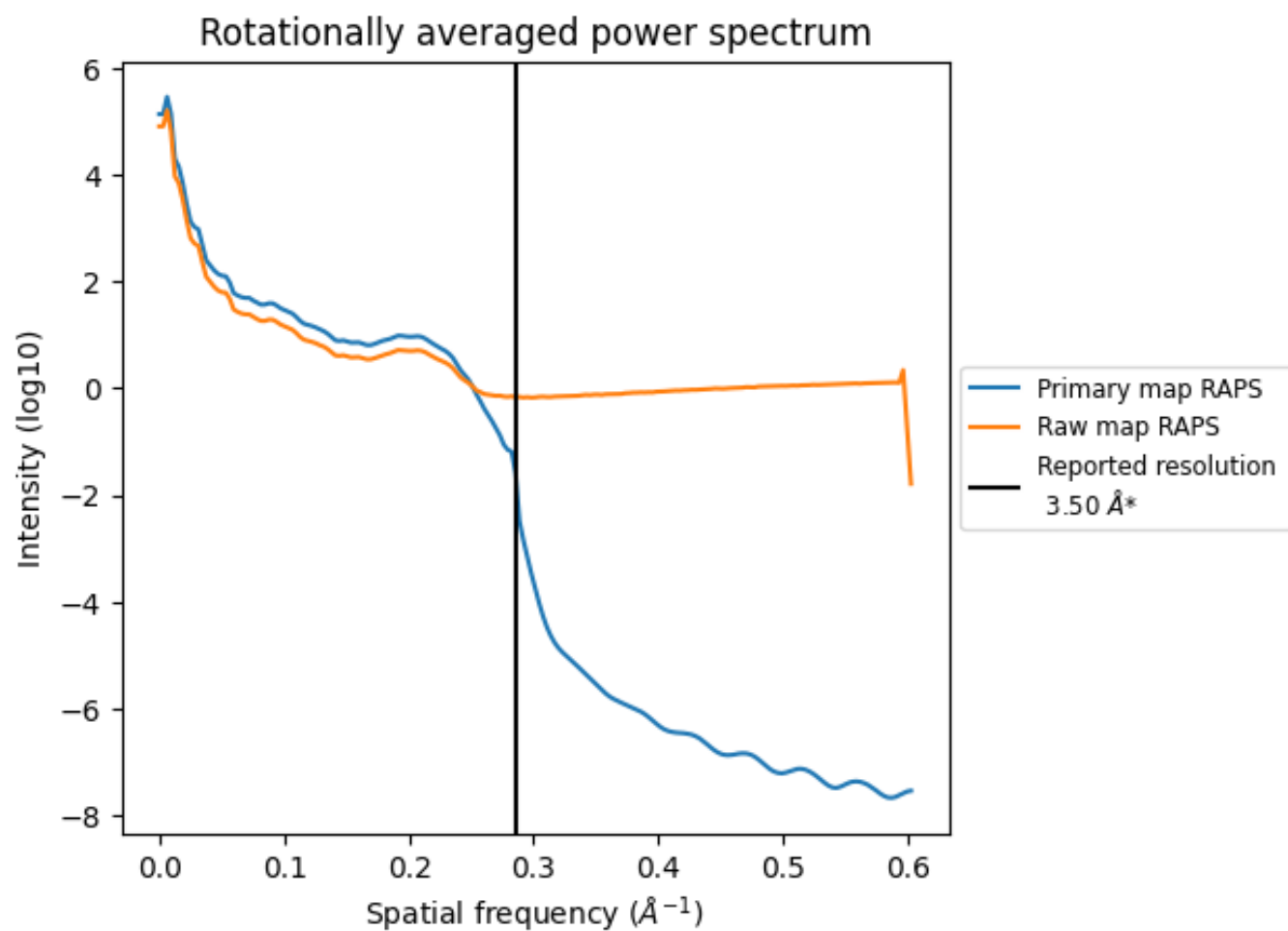
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 174 nm<sup>3</sup>; this corresponds to an approximate mass of 157 kDa.

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

### 7.3 Rotationally averaged power spectrum ⓘ

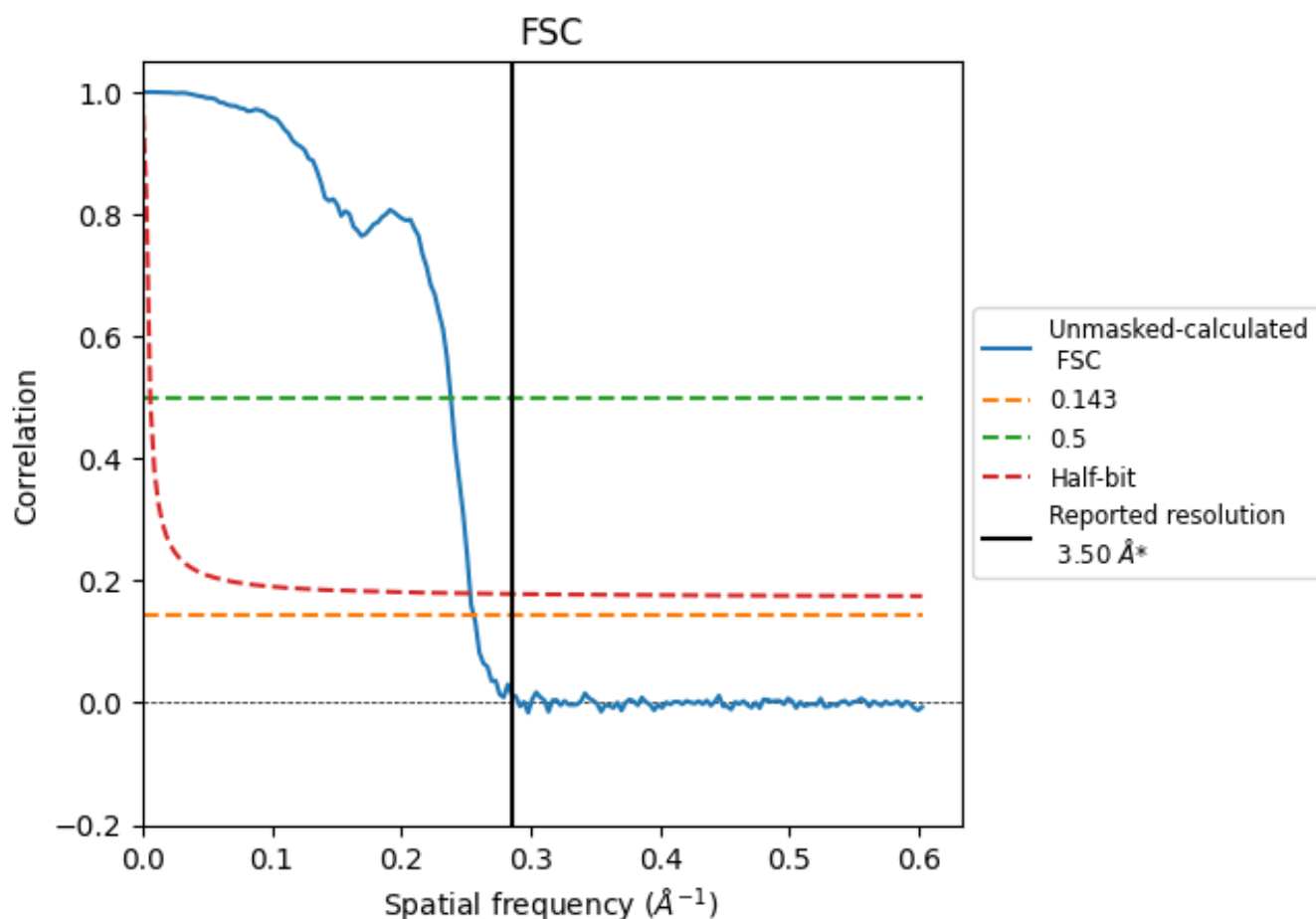


\*Reported resolution corresponds to spatial frequency of 0.286  $\text{\AA}^{-1}$

## 8 Fourier-Shell correlation [i](#)

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

### 8.1 FSC [i](#)



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

## 8.2 Resolution estimates [i](#)

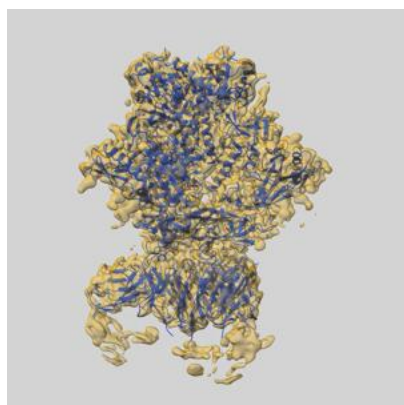
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.90	4.20	3.95

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.90 differs from the reported value 3.5 by more than 10 %

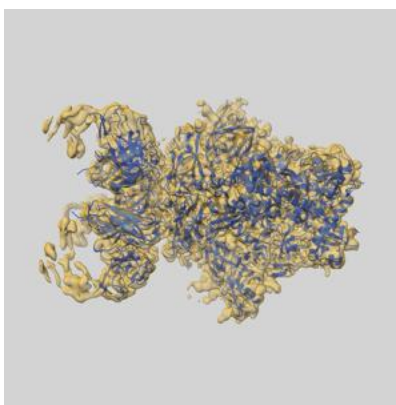
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-70666 and PDB model 9OOM. Per-residue inclusion information can be found in section [3](#) on page [12](#).

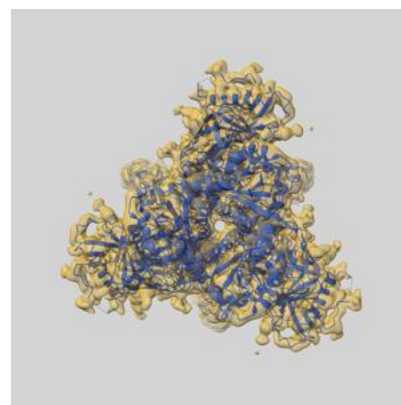
### 9.1 Map-model overlay [i](#)



X



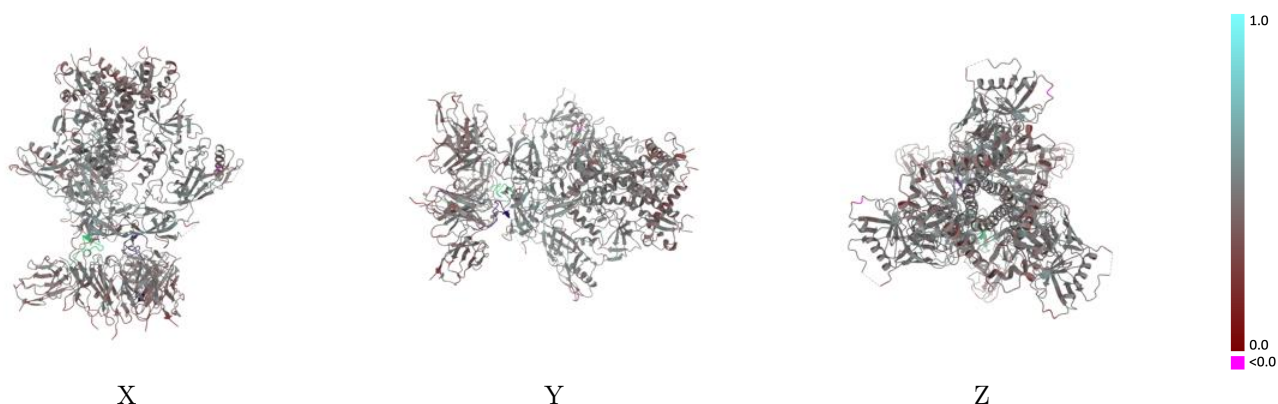
Y



Z

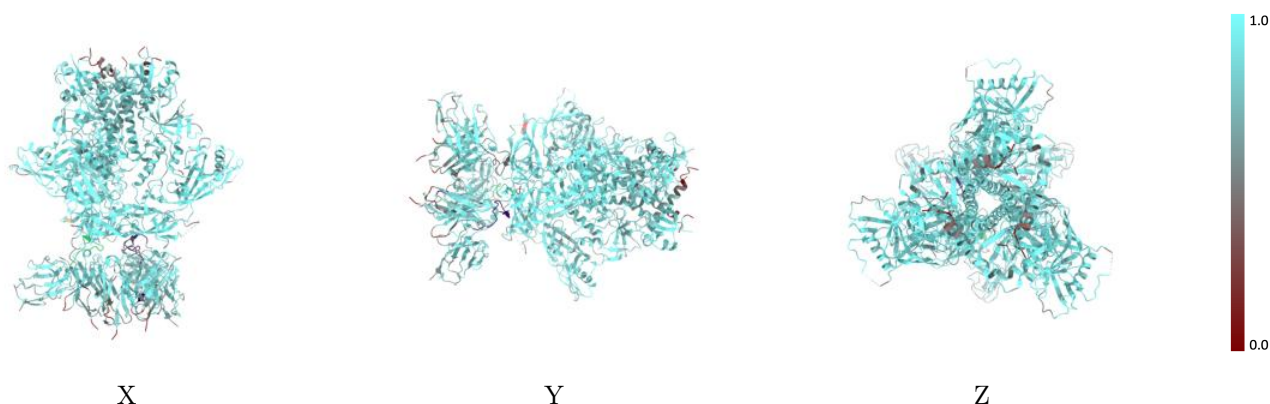
The images above show the 3D surface view of the map at the recommended contour level 0.09 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



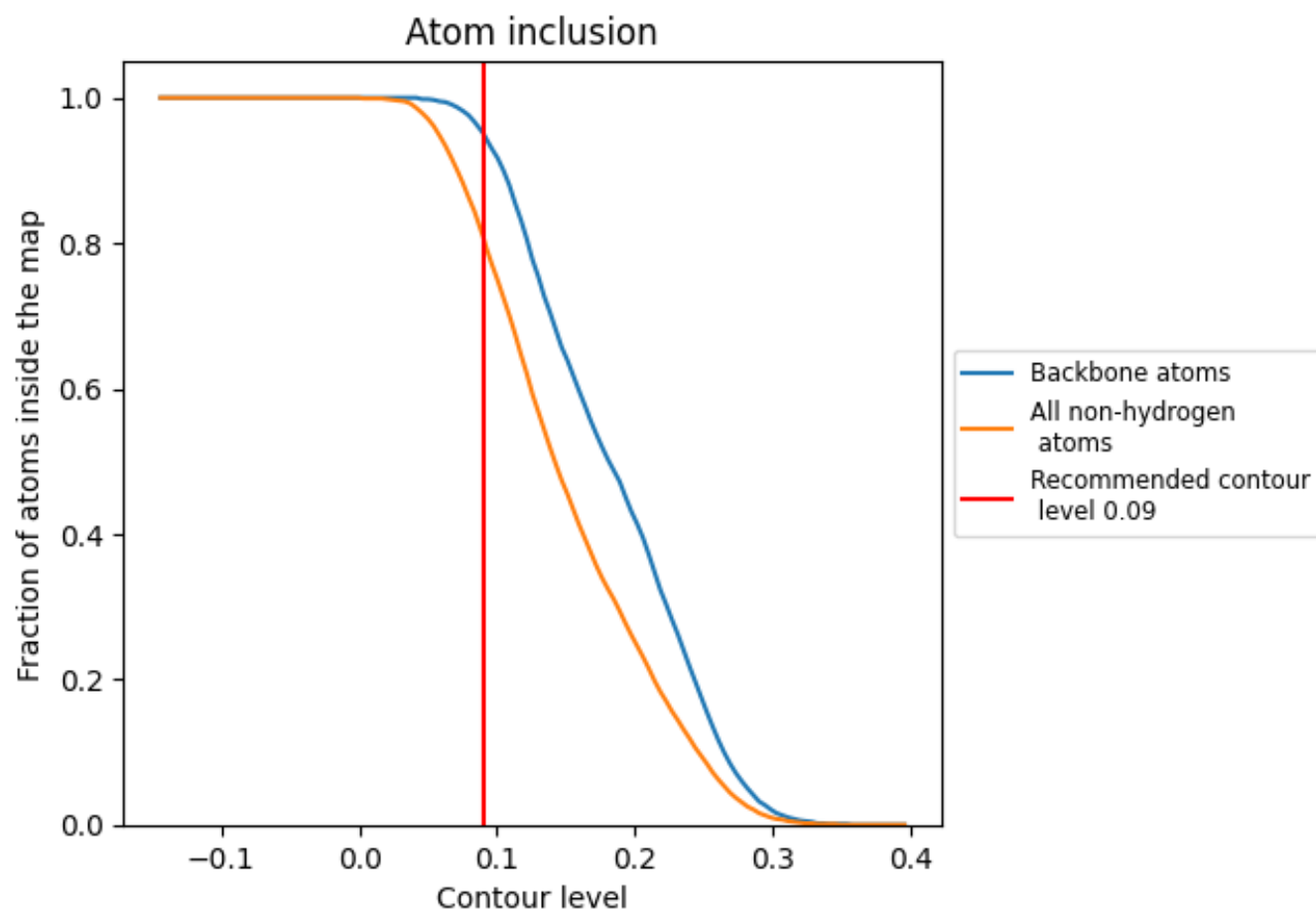
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.09).







































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 81% 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 (0.09) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8080	 0.4450
A	 0.5710	 0.4170
B	 0.7440	 0.4580
C	 0.8530	 0.4760
D	 0.2620	 0.4330
E	 0.7660	 0.4690
F	 0.6790	 0.4630
G	 0.5360	 0.4530
H	 0.7740	 0.4290
I	 0.7770	 0.4290
J	 0.7750	 0.4270
K	 0.3210	 0.3510
L	 0.7580	 0.4130
M	 0.7570	 0.4130
N	 0.7550	 0.4150
O	 0.6430	 0.3950
P	 0.5360	 0.3780
Q	 0.5000	 0.3770
R	 0.5710	 0.3640
S	 0.7500	 0.4130
T	 0.5710	 0.4040
U	 0.7690	 0.4690
V	 0.8030	 0.4790
W	 0.2300	 0.4350
X	 0.8470	 0.4670
Y	 0.6430	 0.4390
Z	 0.5360	 0.4310
a	 0.3210	 0.3120
b	 0.8670	 0.4660
c	 0.7690	 0.4240
d	 0.8670	 0.4660
e	 0.7780	 0.4260
f	 0.8680	 0.4660
g	 0.7690	 0.4230
h	 0.6070	 0.3680



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Chain	Atom inclusion	Q-score
i	 0.5000	 0.3570
j	 0.5360	 0.3340
k	 0.5710	 0.4030
l	 0.6790	 0.3260
m	 0.5710	 0.4210
n	 0.7690	 0.4610
o	 0.8360	 0.4690
p	 0.2460	 0.4270
q	 0.7450	 0.4490
r	 0.6790	 0.4470
s	 0.5000	 0.3950
t	 0.3210	 0.3590
u	 0.6430	 0.3980
v	 0.5360	 0.3640
w	 0.5360	 0.3820
x	 0.5710	 0.3870
y	 0.6790	 0.3770