



# wwPDB EM Validation Summary Report ⓘ

Apr 9, 2025 – 06:39 PM EDT

PDB ID : 9B8F / pdb\_00009b8f  
EMDB ID : EMD-44344  
Title : SARS CoV-2 full-length spike protein with Lys1269Ala and His1271Ala substitutions in the coatomer binding motif, 2RBD-up conformation (SPIKE-AXA)  
Authors : Singh, S.; Hasan, S.S.  
Deposited on : 2024-03-29  
Resolution : 4.15 Å(reported)  
Based on initial model : 7KRQ

This is a wwPDB EM Validation Summary 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.dev117  
Mogul : 2022.3.0, CSD as543be (2022)  
MolProbity : 4.02b-467  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.42

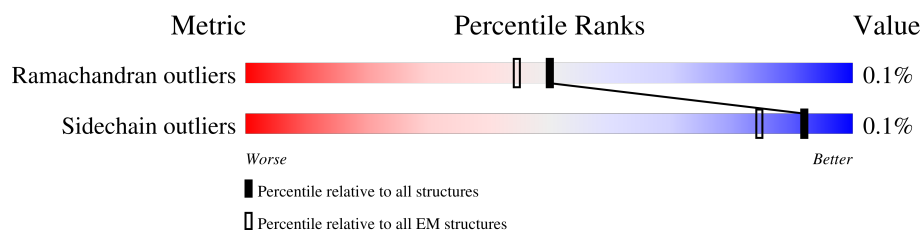
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*




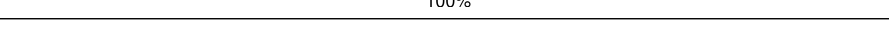
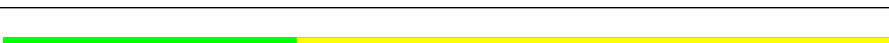

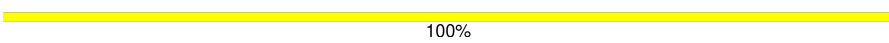


The reported resolution of this entry is 4.15 Å.

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.






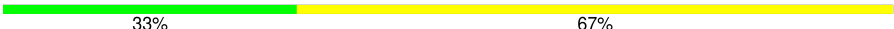
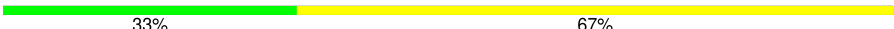
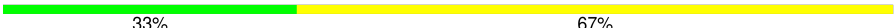
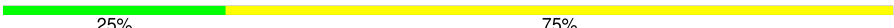

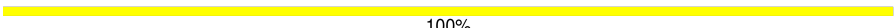
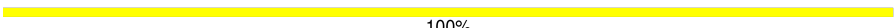
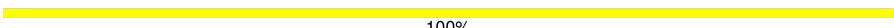
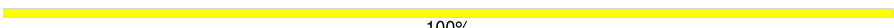
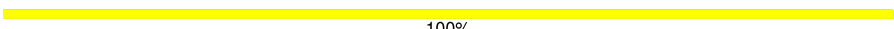
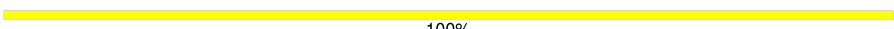
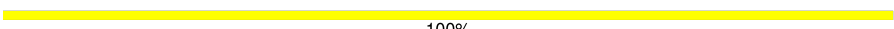
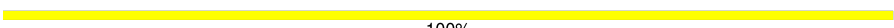

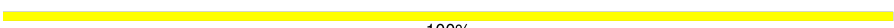

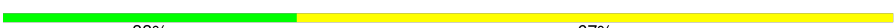
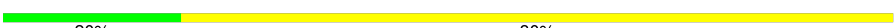


Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	A	1312	
1	B	1312	
1	C	1312	
2	D	3	
2	L	3	
2	M	3	
2	N	3	
2	P	3	
2	R	3	

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Mol	Chain	Length	Quality of chain
2	S	3	 33% 67%
2	U	3	 33% 67%
2	V	3	 33% 67%
2	W	3	 33% 67%
2	X	3	 33% 67%
2	b	3	 33% 67%
3	E	4	 25% 75%
3	F	4	 25% 75%
4	G	2	 100%
4	I	2	 100%
4	K	2	 100%
4	O	2	 100%
4	Q	2	 100%
4	T	2	 100%
4	e	2	 100%
4	f	2	 100%
5	H	4	 50% 50%
6	J	4	 100%
6	Z	4	 25% 75%
7	Y	3	 33% 67%
8	a	5	 20% 80%
8	d	5	 20% 80%
9	c	4	 25% 75%

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 51495 atoms, of which 25203 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 Spike glycoprotein.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	1044	Total	C	H	N	O	S	0	0
			16045	5197	7909	1356	1547	36		
1	B	1076	Total	C	H	N	O	S	0	0
			16504	5360	8106	1399	1601	38		
1	C	1073	Total	C	H	N	O	S	0	0
			16425	5333	8070	1392	1594	36		

There are 141 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	SER	-	insertion	UNP P0DTC2
A	-19	ALA	-	insertion	UNP P0DTC2
A	-18	TRP	-	insertion	UNP P0DTC2
A	-17	SER	-	insertion	UNP P0DTC2
A	-16	HIS	-	insertion	UNP P0DTC2
A	-15	PRO	-	insertion	UNP P0DTC2
A	-14	GLN	-	insertion	UNP P0DTC2
A	-13	PHE	-	insertion	UNP P0DTC2
A	-12	GLU	-	insertion	UNP P0DTC2
A	-11	LYS	-	insertion	UNP P0DTC2
A	-10	GLY	-	insertion	UNP P0DTC2
A	-9	GLY	-	insertion	UNP P0DTC2
A	-8	GLY	-	insertion	UNP P0DTC2
A	-7	SER	-	insertion	UNP P0DTC2
A	-6	GLY	-	insertion	UNP P0DTC2
A	-5	GLY	-	insertion	UNP P0DTC2
A	-4	GLY	-	insertion	UNP P0DTC2
A	-3	SER	-	insertion	UNP P0DTC2
A	-2	GLY	-	insertion	UNP P0DTC2
A	-1	GLY	-	insertion	UNP P0DTC2
A	0	SER	-	insertion	UNP P0DTC2
A	1	SER	-	insertion	UNP P0DTC2
A	2	ALA	-	insertion	UNP P0DTC2
A	3	TRP	-	insertion	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	4	SER	-	insertion	UNP P0DTC2
A	5	HIS	-	insertion	UNP P0DTC2
A	6	PRO	-	insertion	UNP P0DTC2
A	7	GLN	-	insertion	UNP P0DTC2
A	8	PHE	-	insertion	UNP P0DTC2
A	9	GLU	-	insertion	UNP P0DTC2
A	10	LYS	-	insertion	UNP P0DTC2
A	11	SER	-	insertion	UNP P0DTC2
A	12	ALA	-	insertion	UNP P0DTC2
A	13	LEU	-	insertion	UNP P0DTC2
A	14	VAL	-	insertion	UNP P0DTC2
A	15	PRO	-	insertion	UNP P0DTC2
A	16	ARG	-	insertion	UNP P0DTC2
A	17	GLY	-	insertion	UNP P0DTC2
A	18	SER	-	insertion	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	682	GLY	ARG	engineered mutation	UNP P0DTC2
A	683	SER	ARG	engineered mutation	UNP P0DTC2
A	685	SER	ARG	engineered mutation	UNP P0DTC2
A	986	PRO	LYS	engineered mutation	UNP P0DTC2
A	987	PRO	VAL	engineered mutation	UNP P0DTC2
A	1269	ALA	LYS	engineered mutation	UNP P0DTC2
A	1271	ALA	HIS	engineered mutation	UNP P0DTC2
B	-20	SER	-	insertion	UNP P0DTC2
B	-19	ALA	-	insertion	UNP P0DTC2
B	-18	TRP	-	insertion	UNP P0DTC2
B	-17	SER	-	insertion	UNP P0DTC2
B	-16	HIS	-	insertion	UNP P0DTC2
B	-15	PRO	-	insertion	UNP P0DTC2
B	-14	GLN	-	insertion	UNP P0DTC2
B	-13	PHE	-	insertion	UNP P0DTC2
B	-12	GLU	-	insertion	UNP P0DTC2
B	-11	LYS	-	insertion	UNP P0DTC2
B	-10	GLY	-	insertion	UNP P0DTC2
B	-9	GLY	-	insertion	UNP P0DTC2
B	-8	GLY	-	insertion	UNP P0DTC2
B	-7	SER	-	insertion	UNP P0DTC2
B	-6	GLY	-	insertion	UNP P0DTC2
B	-5	GLY	-	insertion	UNP P0DTC2
B	-4	GLY	-	insertion	UNP P0DTC2
B	-3	SER	-	insertion	UNP P0DTC2
B	-2	GLY	-	insertion	UNP P0DTC2

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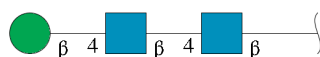
Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	insertion	UNP P0DTC2
B	0	SER	-	insertion	UNP P0DTC2
B	1	SER	-	insertion	UNP P0DTC2
B	2	ALA	-	insertion	UNP P0DTC2
B	3	TRP	-	insertion	UNP P0DTC2
B	4	SER	-	insertion	UNP P0DTC2
B	5	HIS	-	insertion	UNP P0DTC2
B	6	PRO	-	insertion	UNP P0DTC2
B	7	GLN	-	insertion	UNP P0DTC2
B	8	PHE	-	insertion	UNP P0DTC2
B	9	GLU	-	insertion	UNP P0DTC2
B	10	LYS	-	insertion	UNP P0DTC2
B	11	SER	-	insertion	UNP P0DTC2
B	12	ALA	-	insertion	UNP P0DTC2
B	13	LEU	-	insertion	UNP P0DTC2
B	14	VAL	-	insertion	UNP P0DTC2
B	15	PRO	-	insertion	UNP P0DTC2
B	16	ARG	-	insertion	UNP P0DTC2
B	17	GLY	-	insertion	UNP P0DTC2
B	18	SER	-	insertion	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	682	GLY	ARG	engineered mutation	UNP P0DTC2
B	683	SER	ARG	engineered mutation	UNP P0DTC2
B	685	SER	ARG	engineered mutation	UNP P0DTC2
B	986	PRO	LYS	engineered mutation	UNP P0DTC2
B	987	PRO	VAL	engineered mutation	UNP P0DTC2
B	1269	ALA	LYS	engineered mutation	UNP P0DTC2
B	1271	ALA	HIS	engineered mutation	UNP P0DTC2
C	-20	SER	-	insertion	UNP P0DTC2
C	-19	ALA	-	insertion	UNP P0DTC2
C	-18	TRP	-	insertion	UNP P0DTC2
C	-17	SER	-	insertion	UNP P0DTC2
C	-16	HIS	-	insertion	UNP P0DTC2
C	-15	PRO	-	insertion	UNP P0DTC2
C	-14	GLN	-	insertion	UNP P0DTC2
C	-13	PHE	-	insertion	UNP P0DTC2
C	-12	GLU	-	insertion	UNP P0DTC2
C	-11	LYS	-	insertion	UNP P0DTC2
C	-10	GLY	-	insertion	UNP P0DTC2
C	-9	GLY	-	insertion	UNP P0DTC2
C	-8	GLY	-	insertion	UNP P0DTC2
C	-7	SER	-	insertion	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	GLY	-	insertion	UNP P0DTC2
C	-5	GLY	-	insertion	UNP P0DTC2
C	-4	GLY	-	insertion	UNP P0DTC2
C	-3	SER	-	insertion	UNP P0DTC2
C	-2	GLY	-	insertion	UNP P0DTC2
C	-1	GLY	-	insertion	UNP P0DTC2
C	0	SER	-	insertion	UNP P0DTC2
C	1	SER	-	insertion	UNP P0DTC2
C	2	ALA	-	insertion	UNP P0DTC2
C	3	TRP	-	insertion	UNP P0DTC2
C	4	SER	-	insertion	UNP P0DTC2
C	5	HIS	-	insertion	UNP P0DTC2
C	6	PRO	-	insertion	UNP P0DTC2
C	7	GLN	-	insertion	UNP P0DTC2
C	8	PHE	-	insertion	UNP P0DTC2
C	9	GLU	-	insertion	UNP P0DTC2
C	10	LYS	-	insertion	UNP P0DTC2
C	11	SER	-	insertion	UNP P0DTC2
C	12	ALA	-	insertion	UNP P0DTC2
C	13	LEU	-	insertion	UNP P0DTC2
C	14	VAL	-	insertion	UNP P0DTC2
C	15	PRO	-	insertion	UNP P0DTC2
C	16	ARG	-	insertion	UNP P0DTC2
C	17	GLY	-	insertion	UNP P0DTC2
C	18	SER	-	insertion	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	682	GLY	ARG	engineered mutation	UNP P0DTC2
C	683	SER	ARG	engineered mutation	UNP P0DTC2
C	685	SER	ARG	engineered mutation	UNP P0DTC2
C	986	PRO	LYS	engineered mutation	UNP P0DTC2
C	987	PRO	VAL	engineered mutation	UNP P0DTC2
C	1269	ALA	LYS	engineered mutation	UNP P0DTC2
C	1271	ALA	HIS	engineered mutation	UNP P0DTC2

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



Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	3	Total	C	N	O	0	0
			39	22	2	15		
2	L	3	Total	C	H	N	0	0
			73	22	34	2		
2	M	3	Total	C	H	N	0	0
			73	22	34	2		
2	N	3	Total	C	H	N	0	0
			73	22	34	2		
2	P	3	Total	C	H	N	0	0
			73	22	34	2		
2	R	3	Total	C	H	N	0	0
			73	22	34	2		
2	S	3	Total	C	H	N	0	0
			73	22	34	2		
2	U	3	Total	C	H	N	0	0
			73	22	34	2		
2	V	3	Total	C	H	N	0	0
			73	22	34	2		
2	W	3	Total	C	H	N	0	0
			73	22	34	2		
2	X	3	Total	C	H	N	0	0
			73	22	34	2		
2	b	3	Total	C	H	N	0	0
			73	22	34	2		

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



Mol	Chain	Residues	Atoms				AltConf	Trace
3	E	4	Total	C	N	O	0	0
			50	28	2	20		
3	F	4	Total	C	N	O	0	0
			50	28	2	20		

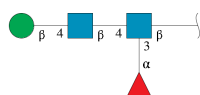
- Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.





Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	2	Total	C	H	N	O	0	0
			53	16	25	2	10		
4	I	2	Total	C	H	N	O	0	0
			53	16	25	2	10		
4	K	2	Total	C	H	N	O	0	0
			53	16	25	2	10		
4	O	2	Total	C	H	N	O	0	0
			53	16	25	2	10		
4	Q	2	Total	C	H	N	O	0	0
			53	16	25	2	10		
4	T	2	Total	C	H	N	O	0	0
			53	16	25	2	10		
4	e	2	Total	C	H	N	O	0	0
			53	16	25	2	10		
4	f	2	Total	C	H	N	O	0	0
			53	16	25	2	10		

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



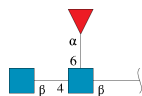
Mol	Chain	Residues	Atoms					AltConf	Trace
5	H	4	Total	C	H	N	O	0	0
			92	28	43	2	19		

- Molecule 6 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.



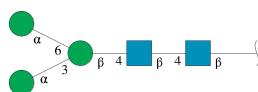
Mol	Chain	Residues	Atoms					AltConf	Trace
6	J	4	Total	C	H	N	O	0	0
			93	28	43	2	20		
6	Z	4	Total	C	H	N	O	0	0
			93	28	43	2	20		

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



Mol	Chain	Residues	Atoms					AltConf	Trace
7	Y	3	Total	C	H	N	O	0	0
			72	22	34	2	14		

- 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	a	5	Total	C	H	N	O	0	0
			113	34	52	2	25		
8	d	5	Total	C	H	N	O	0	0
			113	34	52	2	25		

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



Mol	Chain	Residues	Atoms					AltConf	Trace
9	c	4	Total	C	H	N	O	0	0
			93	28	43	2	20		

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



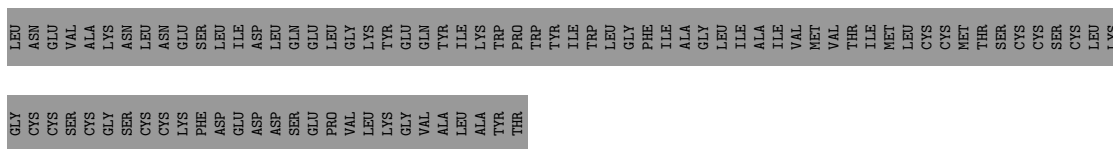
Mol	Chain	Residues	Atoms					AltConf
10	A	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	A	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	A	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	A	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	A	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	B	1	Total	C	H	N	O	0
			27	8	13	1	5	

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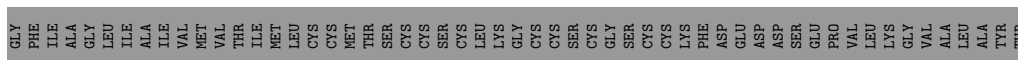
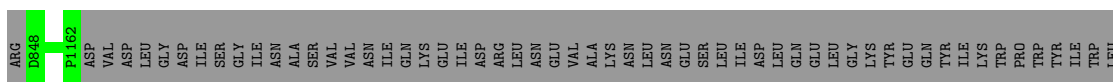
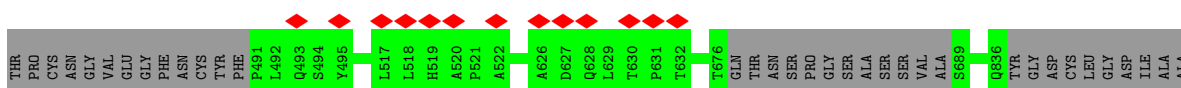
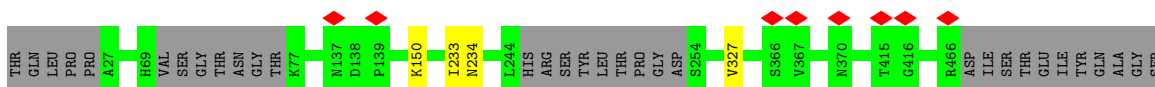
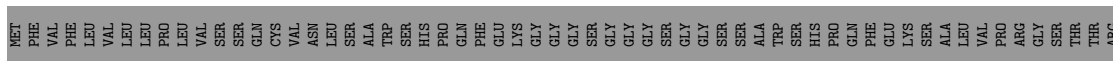
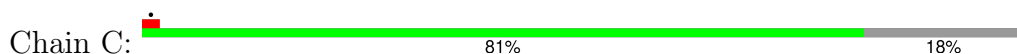
*Continued from previous page...*

Mol	Chain	Residues	Atoms					AltConf
10	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
10	C	1	Total	C	H	N	O	0
			27	8	13	1	5	





- Molecule 1: Spike glycoprotein



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




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



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

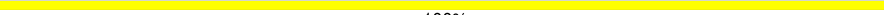


- Molecule 2: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  33% 67%

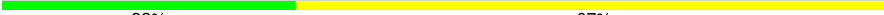
NAG1  
NAG2  
BMA3

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

Chain P:  100%


NAG1  
NAG2  
BMA3

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

Chain R:  33% 67%


NAG1  
NAG2  
BMA3

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

Chain S:  33% 67%


NAG1  
NAG2  
BMA3

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

Chain U:  33% 67%

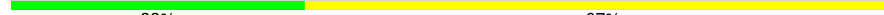
NAG1  
NAG2  
BMA3

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

Chain V:  33% 67%

NAG1  
NAG2  
BMA3

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

Chain W:  33% 67%

NAG1  
NAG2  
BMA3

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

Chain X:  33% 67%

MAG1  
MAG2  
BMA3

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

Chain b:  33% 67%

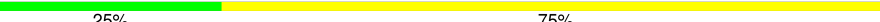
MAG1  
MAG2  
BMA3

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

Chain E:  25% 75%

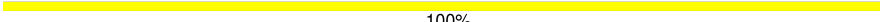
MAG1  
MAG2  
BMA3  
MAN4

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

Chain F:  25% 75%

MAG1  
MAG2  
BMA3  
MAN4

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

Chain G:  100%

MAG1  
MAG2

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

Chain I:  100%

MAG1  
MAG2

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

Chain K:  100%



MAG1  
MAG2

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

Chain O:  100%MAG1  
MAG2

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

Chain Q:  100%MAG1  
MAG2

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

Chain T:  100%MAG1  
MAG2

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

Chain e:  100%MAG1  
MAG2

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

Chain f:  100%MAG1  
MAG2

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



Chain H:  50% 50%MAG1  
MAG2  
EMJ3  
FUC4

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

Chain J:  100%

MAG1  
MAG2  
BMA3  
MAN4

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

Chain Z:  25%  75%

MAG1  
MAG2  
BMA3  
MAN4

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

Chain Y:  33%  67%

MAG1  
MAG2  
FUC3

- 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 a:  20%  80%

MAG1  
MAG2  
BMA3  
MAN4  
MAN5

- 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:  20%  80%

MAG1  
MAG2  
BMA3  
MAN4  
MAN5

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

Chain c:  25%  75%

MAG1  
MAG2  
BMA3  
MAN4

## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	59137	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	42.19	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	105000	Depositor
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	0.425	Depositor
Minimum map value	-0.135	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.018	Depositor
Recommended contour level	0.0607	Depositor
Map size (Å)	332.112, 332.112, 332.112	wwPDB
Map dimensions	272, 272, 272	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.221, 1.221, 1.221	Depositor

## 5 Model quality

### 5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.27	0/8318	0.49	0/11312
1	B	0.27	0/8594	0.49	1/11694 (0.0%)
1	C	0.26	0/8549	0.48	0/11638
All	All	0.27	0/25461	0.49	1/34644 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
All	All	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	165	ASN	N-CA-CB	5.57	120.63	110.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1133	VAL	Peptide
1	B	165	ASN	Peptide
1	C	233	ILE	Peptide

## 5.2 Too-close contacts ⓘ

Due to software issues we are unable to calculate clashes - this section is therefore empty.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1028/1312 (78%)	950 (92%)	78 (8%)	0	100	100
1	B	1064/1312 (81%)	968 (91%)	95 (9%)	1 (0%)	48	82
1	C	1061/1312 (81%)	941 (89%)	117 (11%)	3 (0%)	37	72
All	All	3153/3936 (80%)	2859 (91%)	290 (9%)	4 (0%)	50	82

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	150	LYS
1	C	327	VAL
1	C	234	ASN
1	B	332	ILE

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	905/1135 (80%)	903 (100%)	2 (0%)	92	94
1	B	931/1135 (82%)	929 (100%)	2 (0%)	92	94
1	C	925/1135 (82%)	925 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2761/3405 (81%)	2757 (100%)	4 (0%)	92 95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	LYS
1	A	466	ARG
1	B	328	ARG
1	B	1154	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	928	ASN
1	C	544	ASN
1	C	1011	GLN
1	C	925	ASN
1	A	824	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

89 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	D	1	1,2	14,14,15	1.99	4 (28%)	17,19,21	1.10	2 (11%)
2	NAG	D	2	2	14,14,15	2.05	4 (28%)	17,19,21	1.43	2 (11%)
2	BMA	D	3	2	11,11,12	0.90	1 (9%)	15,15,17	0.92	1 (6%)
3	NAG	E	1	3,1	14,14,15	2.05	5 (35%)	17,19,21	1.03	2 (11%)
3	NAG	E	2	3	14,14,15	2.01	3 (21%)	17,19,21	1.14	2 (11%)
3	BMA	E	3	3	11,11,12	0.46	0	15,15,17	0.71	0
3	MAN	E	4	3	11,11,12	0.86	0	15,15,17	1.49	2 (13%)
3	NAG	F	1	3,1	14,14,15	2.30	4 (28%)	17,19,21	1.96	4 (23%)
3	NAG	F	2	3	14,14,15	2.14	4 (28%)	17,19,21	1.45	3 (17%)
3	BMA	F	3	3	11,11,12	0.66	0	15,15,17	0.84	0
3	MAN	F	4	3	11,11,12	0.89	1 (9%)	15,15,17	1.57	2 (13%)
4	NAG	G	1	4,1	14,14,15	2.42	5 (35%)	17,19,21	2.58	8 (47%)
4	NAG	G	2	4	14,14,15	2.02	4 (28%)	17,19,21	1.25	1 (5%)
5	NAG	H	1	1,5	14,14,15	2.14	4 (28%)	17,19,21	1.56	3 (17%)
5	NAG	H	2	5	14,14,15	1.95	3 (21%)	17,19,21	1.97	5 (29%)
5	BMA	H	3	5	11,11,12	0.54	0	15,15,17	0.66	0
5	FUC	H	4	5	10,10,11	0.76	0	14,14,16	0.65	0
4	NAG	I	1	4,1	14,14,15	2.09	5 (35%)	17,19,21	1.32	1 (5%)
4	NAG	I	2	4	14,14,15	2.09	4 (28%)	17,19,21	1.12	1 (5%)
6	NAG	J	1	1,6	14,14,15	2.07	5 (35%)	17,19,21	1.28	2 (11%)
6	NAG	J	2	6	14,14,15	2.03	5 (35%)	17,19,21	1.18	2 (11%)
6	BMA	J	3	6	11,11,12	1.12	1 (9%)	15,15,17	0.87	0
6	MAN	J	4	6	11,11,12	1.28	3 (27%)	15,15,17	2.39	3 (20%)
4	NAG	K	1	4,1	14,14,15	2.07	4 (28%)	17,19,21	1.71	4 (23%)
4	NAG	K	2	4	14,14,15	1.94	4 (28%)	17,19,21	1.11	2 (11%)
2	NAG	L	1	1,2	14,14,15	2.16	4 (28%)	17,19,21	1.48	2 (11%)
2	NAG	L	2	2	14,14,15	1.95	4 (28%)	17,19,21	1.30	1 (5%)
2	BMA	L	3	2	11,11,12	0.95	1 (9%)	15,15,17	0.84	0
2	NAG	M	1	1,2	14,14,15	2.28	6 (42%)	17,19,21	2.38	8 (47%)
2	NAG	M	2	2	14,14,15	1.96	4 (28%)	17,19,21	1.57	3 (17%)
2	BMA	M	3	2	11,11,12	0.46	0	15,15,17	0.74	0
2	NAG	N	1	1,2	14,14,15	2.05	4 (28%)	17,19,21	1.55	3 (17%)
2	NAG	N	2	2	14,14,15	1.92	4 (28%)	17,19,21	1.23	1 (5%)
2	BMA	N	3	2	11,11,12	0.68	0	15,15,17	0.72	0
4	NAG	O	1	4	14,14,15	1.95	5 (35%)	17,19,21	1.33	2 (11%)
4	NAG	O	2	4	14,14,15	1.99	4 (28%)	17,19,21	1.09	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	P	1	1,2	14,14,15	2.03	4 (28%)	17,19,21	1.39	3 (17%)
2	NAG	P	2	2	14,14,15	2.01	4 (28%)	17,19,21	1.41	3 (17%)
2	BMA	P	3	2	11,11,12	0.64	0	15,15,17	0.86	1 (6%)
4	NAG	Q	1	4,1	14,14,15	2.05	4 (28%)	17,19,21	1.21	2 (11%)
4	NAG	Q	2	4	14,14,15	2.07	4 (28%)	17,19,21	1.44	3 (17%)
2	NAG	R	1	1,2	14,14,15	2.07	4 (28%)	17,19,21	1.28	2 (11%)
2	NAG	R	2	2	14,14,15	2.26	5 (35%)	17,19,21	2.22	5 (29%)
2	BMA	R	3	2	11,11,12	0.62	0	15,15,17	0.77	0
2	NAG	S	1	1,2	14,14,15	2.09	4 (28%)	17,19,21	2.04	6 (35%)
2	NAG	S	2	2	14,14,15	1.98	4 (28%)	17,19,21	1.31	2 (11%)
2	BMA	S	3	2	11,11,12	0.55	0	15,15,17	0.75	0
4	NAG	T	1	4,1	14,14,15	1.98	3 (21%)	17,19,21	1.28	1 (5%)
4	NAG	T	2	4	14,14,15	1.97	4 (28%)	17,19,21	1.29	2 (11%)
2	NAG	U	1	1,2	14,14,15	2.05	4 (28%)	17,19,21	1.60	4 (23%)
2	NAG	U	2	2	14,14,15	2.05	4 (28%)	17,19,21	1.26	2 (11%)
2	BMA	U	3	2	11,11,12	0.65	0	15,15,17	0.69	0
2	NAG	V	1	1,2	14,14,15	1.96	3 (21%)	17,19,21	2.14	4 (23%)
2	NAG	V	2	2	14,14,15	1.94	3 (21%)	17,19,21	1.22	2 (11%)
2	BMA	V	3	2	11,11,12	0.72	0	15,15,17	0.69	0
2	NAG	W	1	1,2	14,14,15	1.89	3 (21%)	17,19,21	2.13	7 (41%)
2	NAG	W	2	2	14,14,15	1.92	3 (21%)	17,19,21	1.36	3 (17%)
2	BMA	W	3	2	11,11,12	0.57	0	15,15,17	0.68	0
2	NAG	X	1	1,2	14,14,15	2.00	4 (28%)	17,19,21	1.35	2 (11%)
2	NAG	X	2	2	14,14,15	1.98	4 (28%)	17,19,21	1.47	2 (11%)
2	BMA	X	3	2	11,11,12	0.59	0	15,15,17	0.68	0
7	NAG	Y	1	7	14,14,15	1.97	5 (35%)	17,19,21	1.31	2 (11%)
7	NAG	Y	2	7	14,14,15	2.06	3 (21%)	17,19,21	1.42	3 (17%)
7	FUC	Y	3	7	10,10,11	0.60	0	14,14,16	0.73	0
6	NAG	Z	1	1,6	14,14,15	1.94	5 (35%)	17,19,21	1.26	2 (11%)
6	NAG	Z	2	6	14,14,15	1.95	3 (21%)	17,19,21	1.00	1 (5%)
6	BMA	Z	3	6	11,11,12	0.53	0	15,15,17	0.86	0
6	MAN	Z	4	6	11,11,12	0.59	0	15,15,17	0.99	2 (13%)
8	NAG	a	1	8,1	14,14,15	2.05	5 (35%)	17,19,21	1.16	2 (11%)
8	NAG	a	2	8	14,14,15	2.15	4 (28%)	17,19,21	1.62	4 (23%)
8	BMA	a	3	8	11,11,12	0.58	0	15,15,17	0.71	0
8	MAN	a	4	8	11,11,12	0.76	0	15,15,17	1.41	3 (20%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
8	MAN	a	5	8	11,11,12	0.64	0	15,15,17	1.32	1 (6%)
2	NAG	b	1	2	14,14,15	2.04	5 (35%)	17,19,21	1.40	3 (17%)
2	NAG	b	2	2	14,14,15	2.08	4 (28%)	17,19,21	1.20	2 (11%)
2	BMA	b	3	2	11,11,12	0.63	0	15,15,17	0.68	0
9	NAG	c	1	1,9	14,14,15	1.83	5 (35%)	17,19,21	2.29	6 (35%)
9	NAG	c	2	9	14,14,15	1.90	3 (21%)	17,19,21	2.06	4 (23%)
9	BMA	c	3	9	11,11,12	0.82	0	15,15,17	0.81	0
9	MAN	c	4	9	11,11,12	0.74	0	15,15,17	1.16	2 (13%)
8	NAG	d	1	8	14,14,15	2.02	4 (28%)	17,19,21	1.46	2 (11%)
8	NAG	d	2	8	14,14,15	1.94	3 (21%)	17,19,21	1.20	2 (11%)
8	BMA	d	3	8	11,11,12	0.81	0	15,15,17	0.79	0
8	MAN	d	4	8	11,11,12	0.75	0	15,15,17	1.18	2 (13%)
8	MAN	d	5	8	11,11,12	0.67	0	15,15,17	0.84	1 (6%)
4	NAG	e	1	4,1	14,14,15	1.94	4 (28%)	17,19,21	1.29	2 (11%)
4	NAG	e	2	4	14,14,15	2.00	4 (28%)	17,19,21	1.08	1 (5%)
4	NAG	f	1	4,1	14,14,15	2.08	5 (35%)	17,19,21	1.25	2 (11%)
4	NAG	f	2	4	14,14,15	2.06	4 (28%)	17,19,21	1.32	2 (11%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	D	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	D	2	2	-	1/6/23/26	0/1/1/1
2	BMA	D	3	2	-	0/2/19/22	0/1/1/1
3	NAG	E	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	BMA	E	3	3	-	2/2/19/22	0/1/1/1
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
3	NAG	F	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	2/2/19/22	0/1/1/1
3	MAN	F	4	3	-	0/2/19/22	0/1/1/1
4	NAG	G	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	3/6/23/26	0/1/1/1
5	NAG	H	1	1,5	-	3/6/23/26	0/1/1/1

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

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

The worst 5 of 243 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	1	NAG	O5-C1	5.60	1.53	1.43
2	R	2	NAG	O5-C1	5.34	1.52	1.43
3	F	2	NAG	O5-C1	4.86	1.51	1.43
2	S	1	NAG	O5-C1	4.73	1.51	1.43
3	F	1	NAG	C7-N2	4.71	1.49	1.34

The worst 5 of 181 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	4	MAN	C1-O5-C5	8.12	123.07	112.19
4	G	1	NAG	C1-O5-C5	6.81	121.31	112.19
2	V	1	NAG	C1-O5-C5	-6.64	103.28	112.19
3	F	1	NAG	C1-O5-C5	-6.36	103.67	112.19
2	R	2	NAG	C1-O5-C5	5.69	119.81	112.19

There are no chirality outliers.

5 of 136 torsion outliers are listed below:

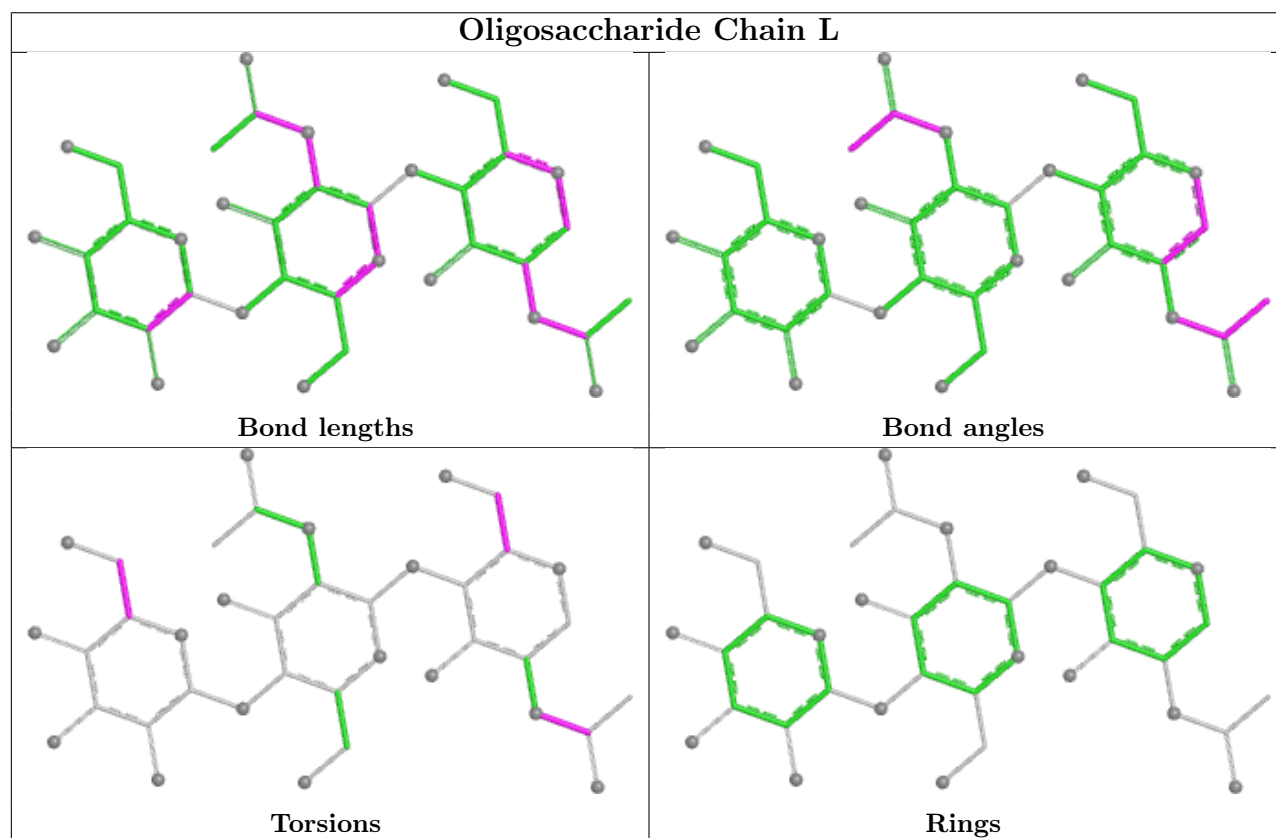
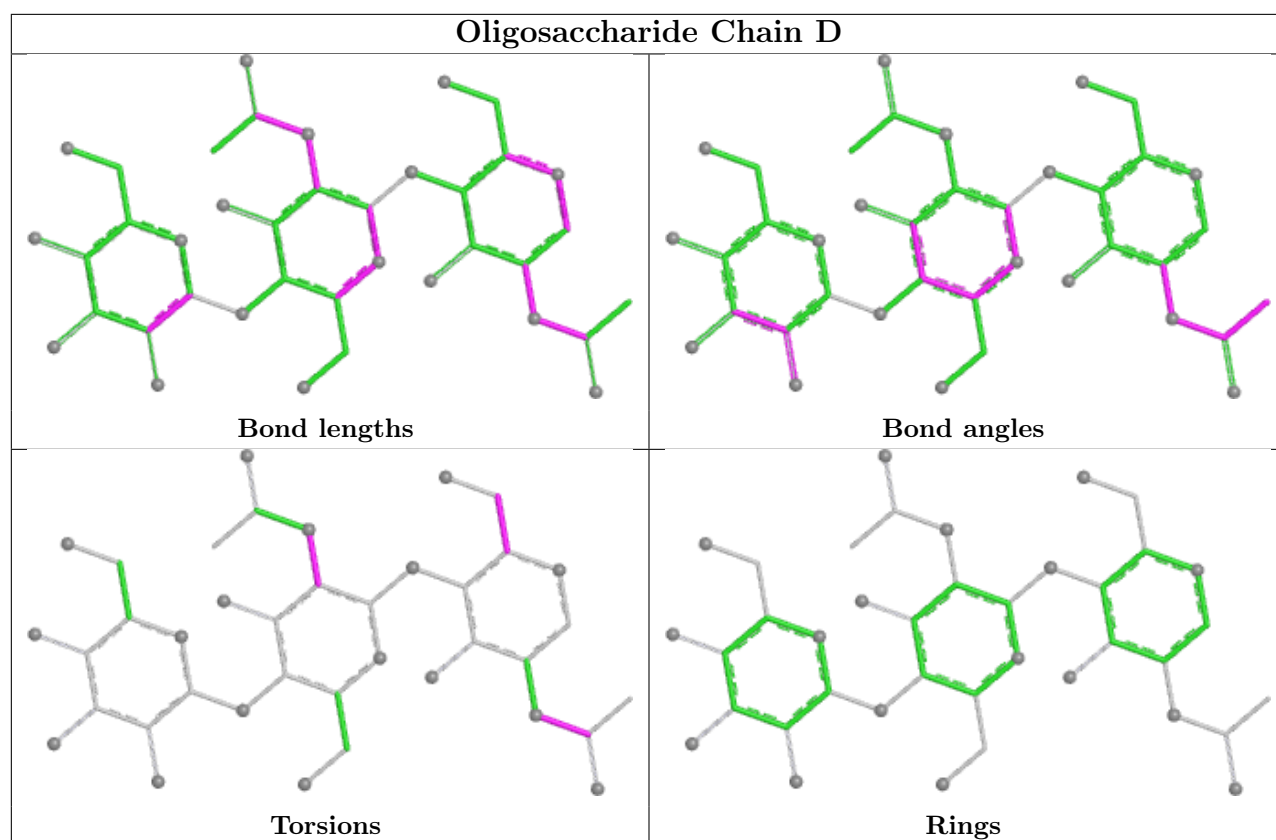
Mol	Chain	Res	Type	Atoms
2	D	2	NAG	C3-C2-N2-C7
2	N	1	NAG	C1-C2-N2-C7
2	V	1	NAG	C1-C2-N2-C7
4	G	2	NAG	C3-C2-N2-C7
4	e	1	NAG	C1-C2-N2-C7

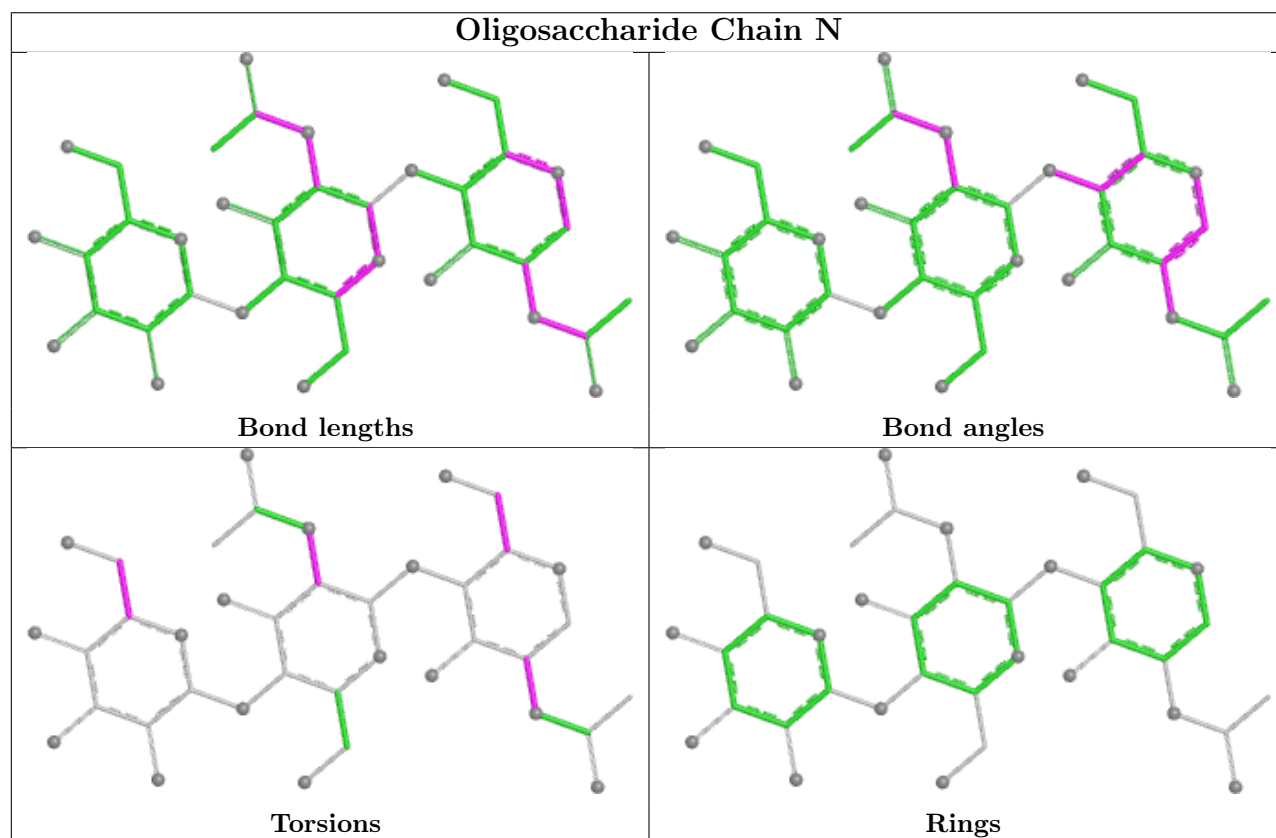
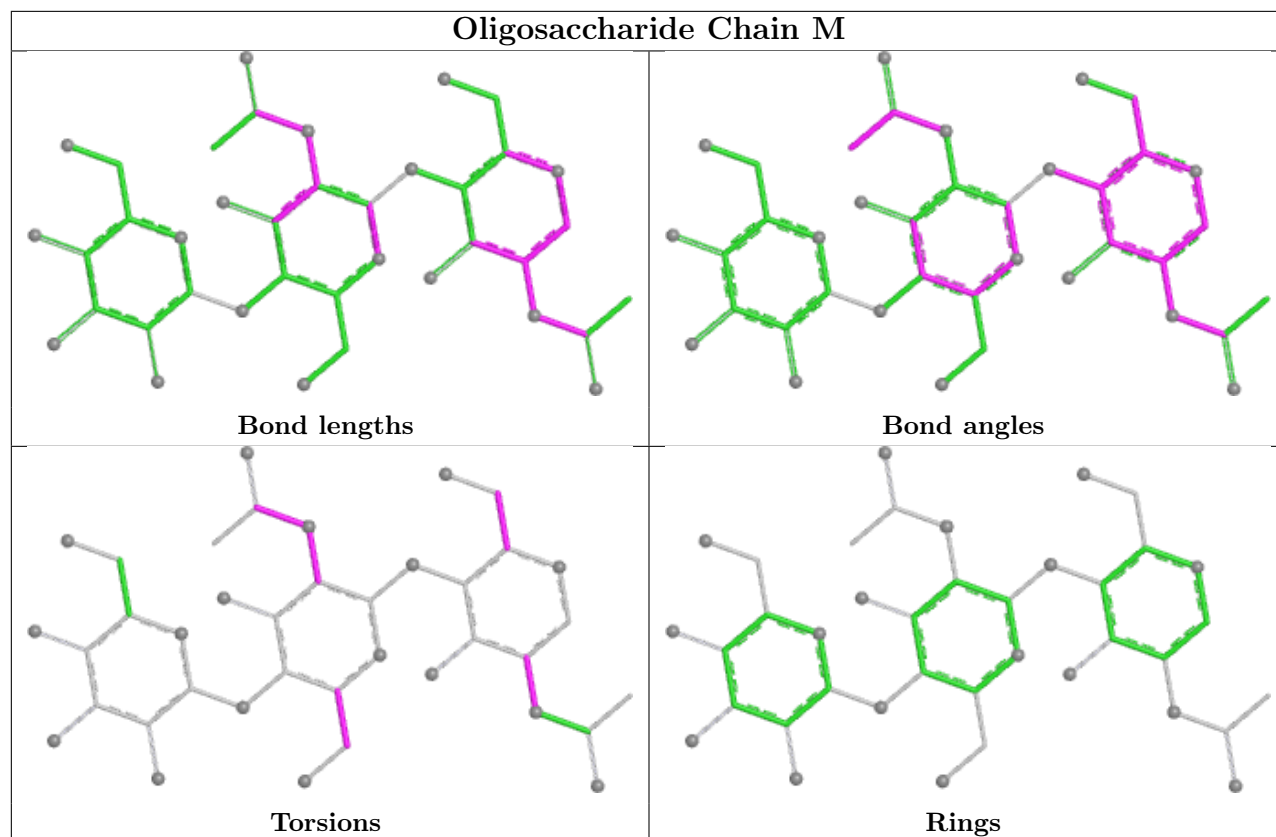
All (2) ring outliers are listed below:

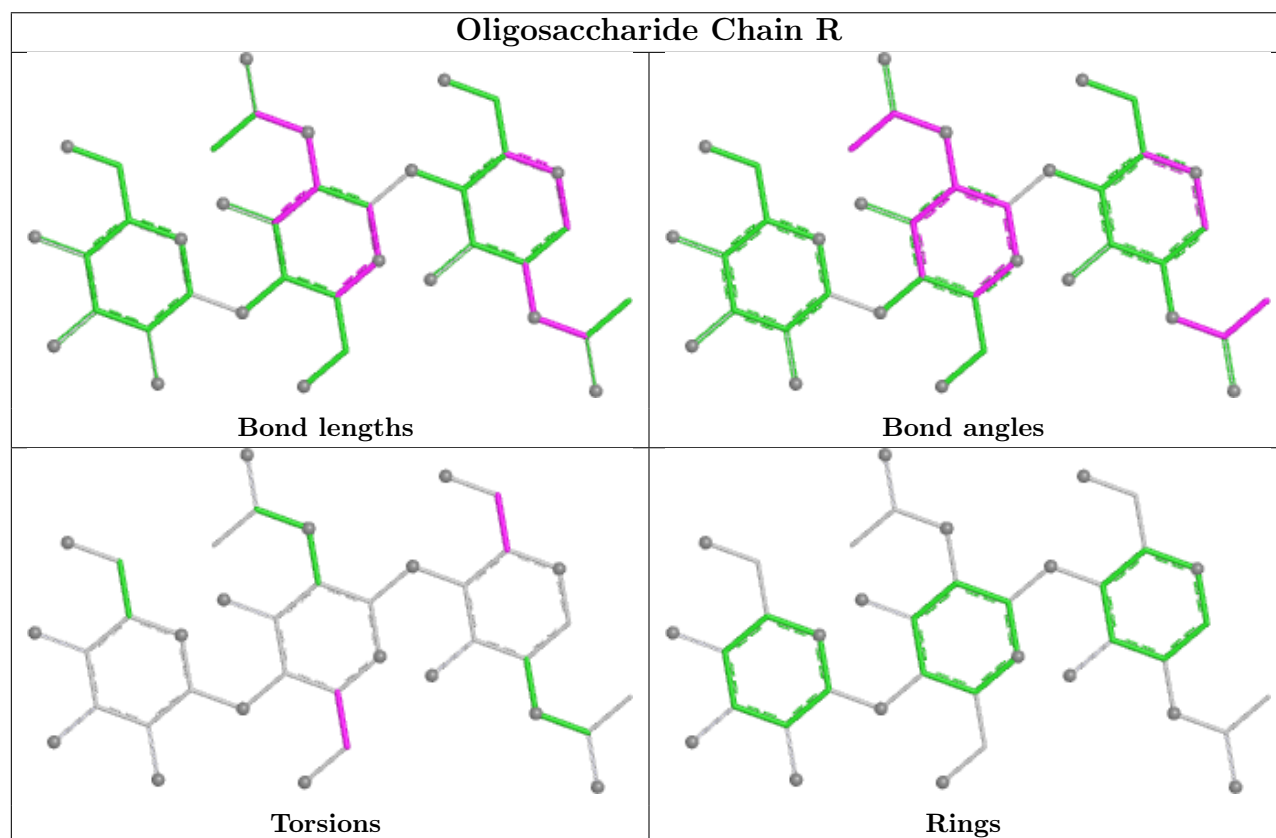
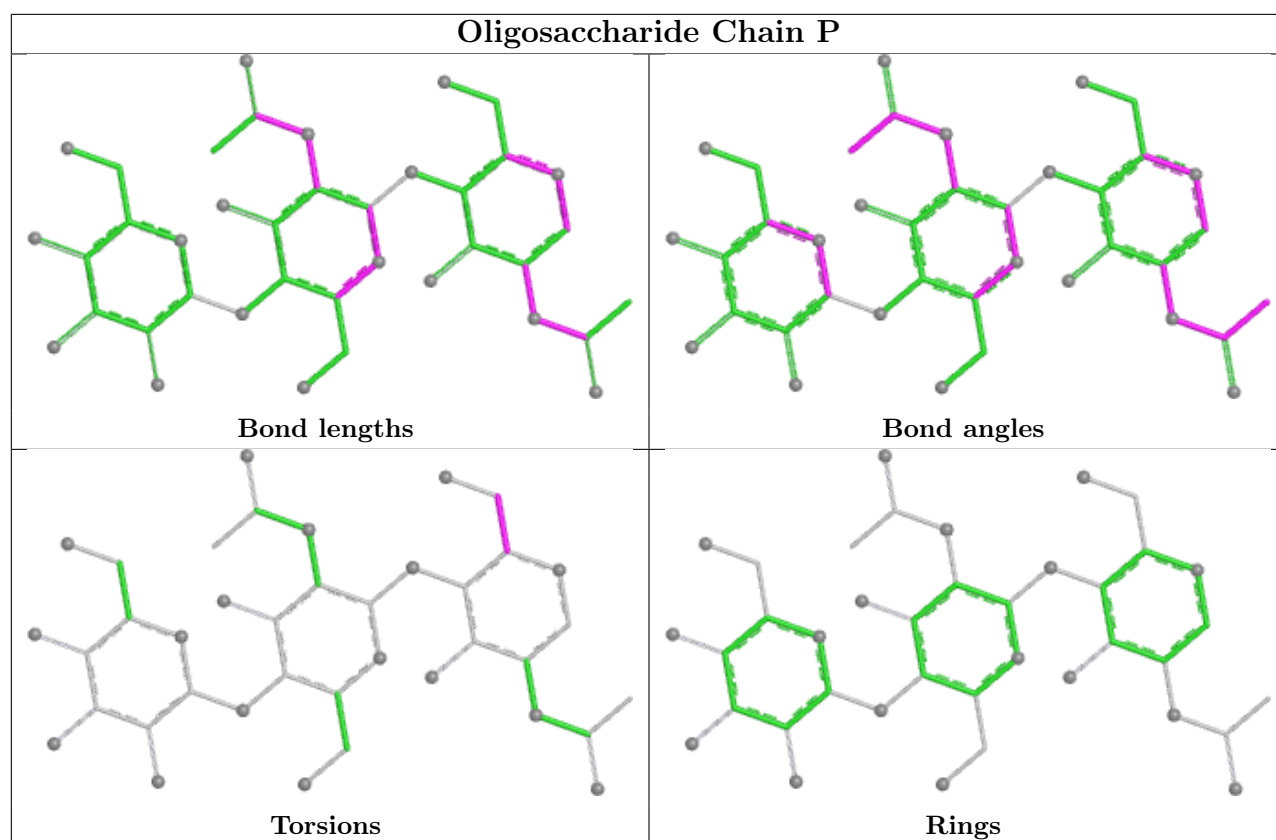
Mol	Chain	Res	Type	Atoms
8	a	3	BMA	C1-C2-C3-C4-C5-O5
8	a	5	MAN	C1-C2-C3-C4-C5-O5

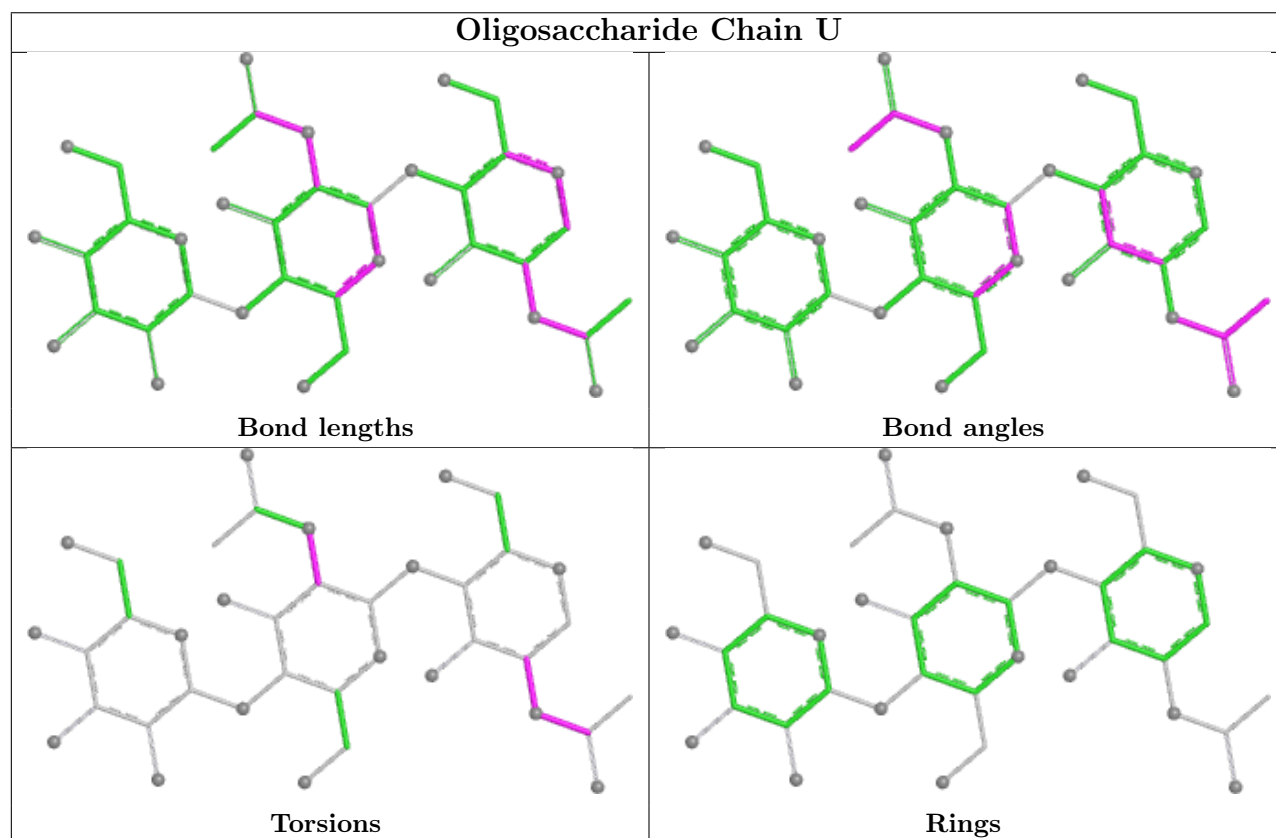
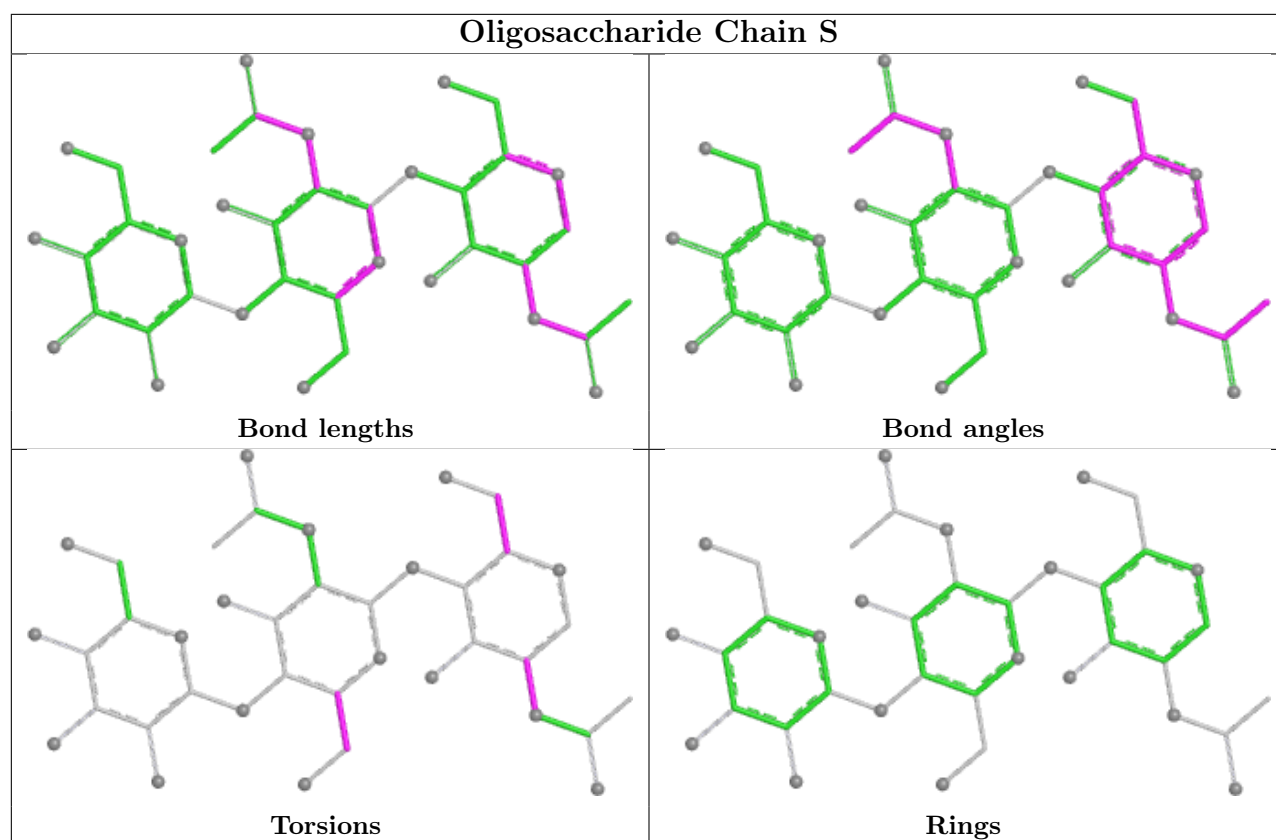
No monomer is involved in short contacts.

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

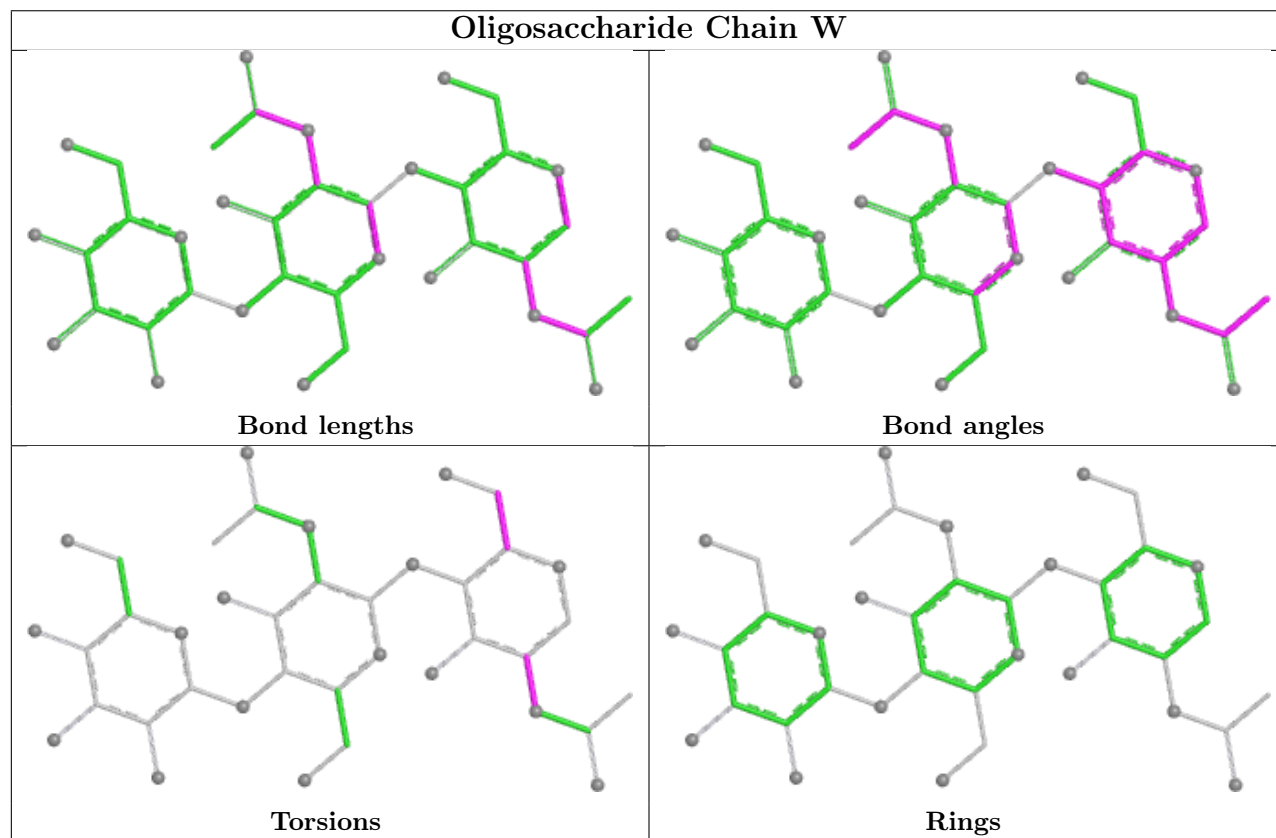
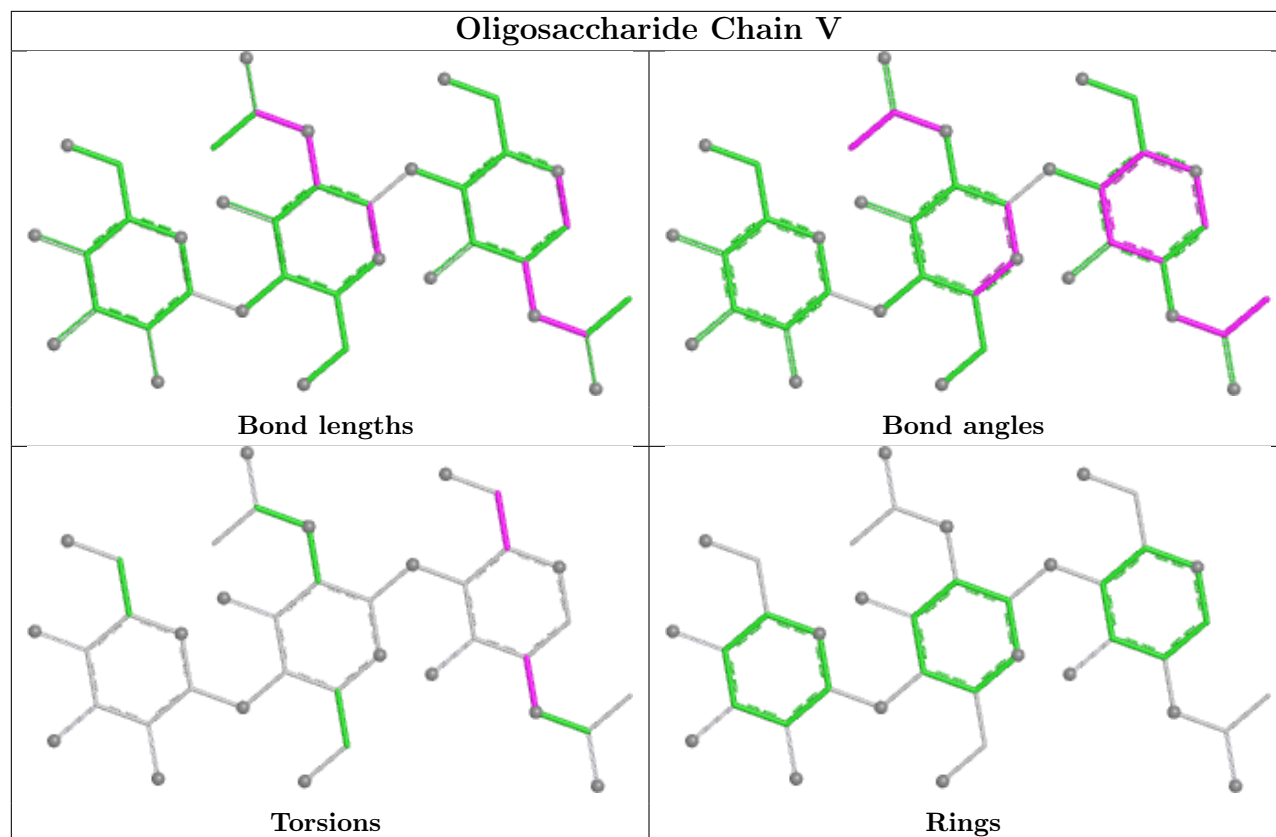


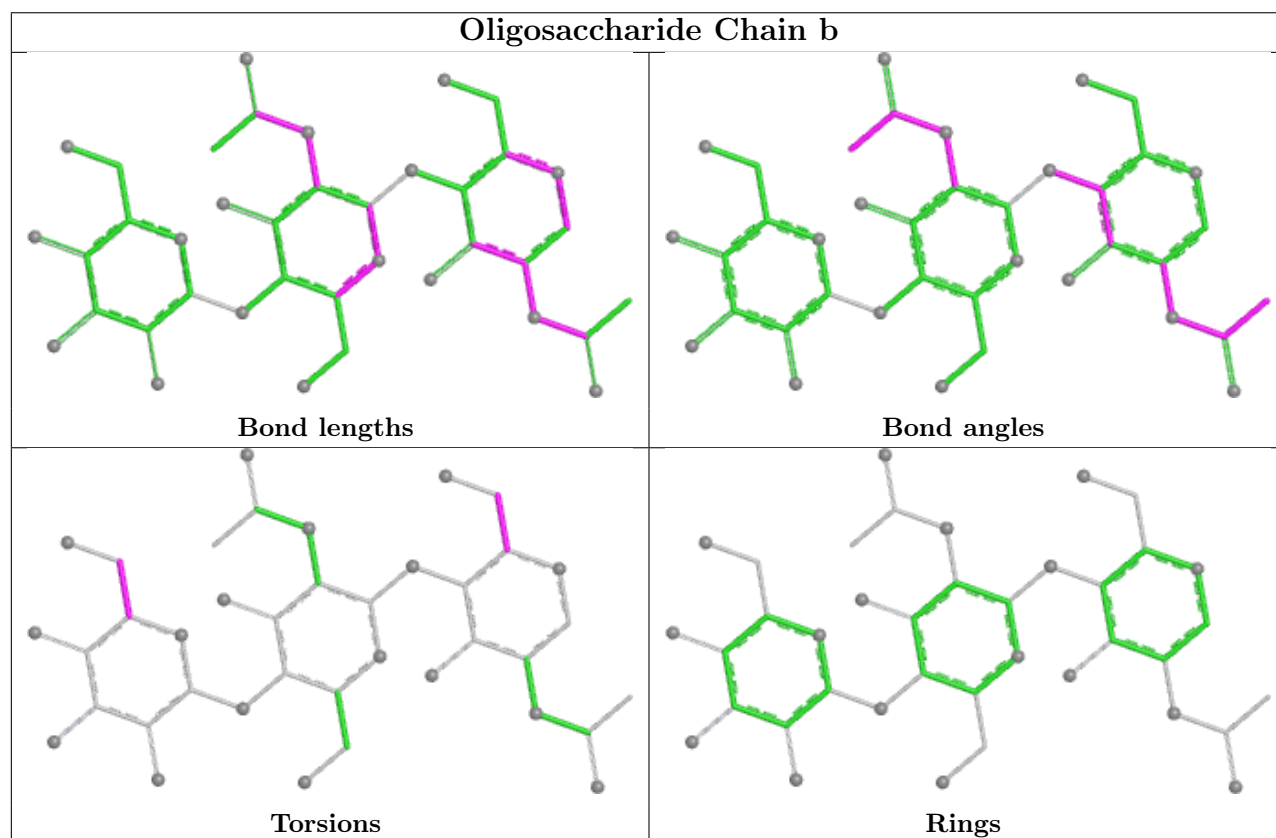
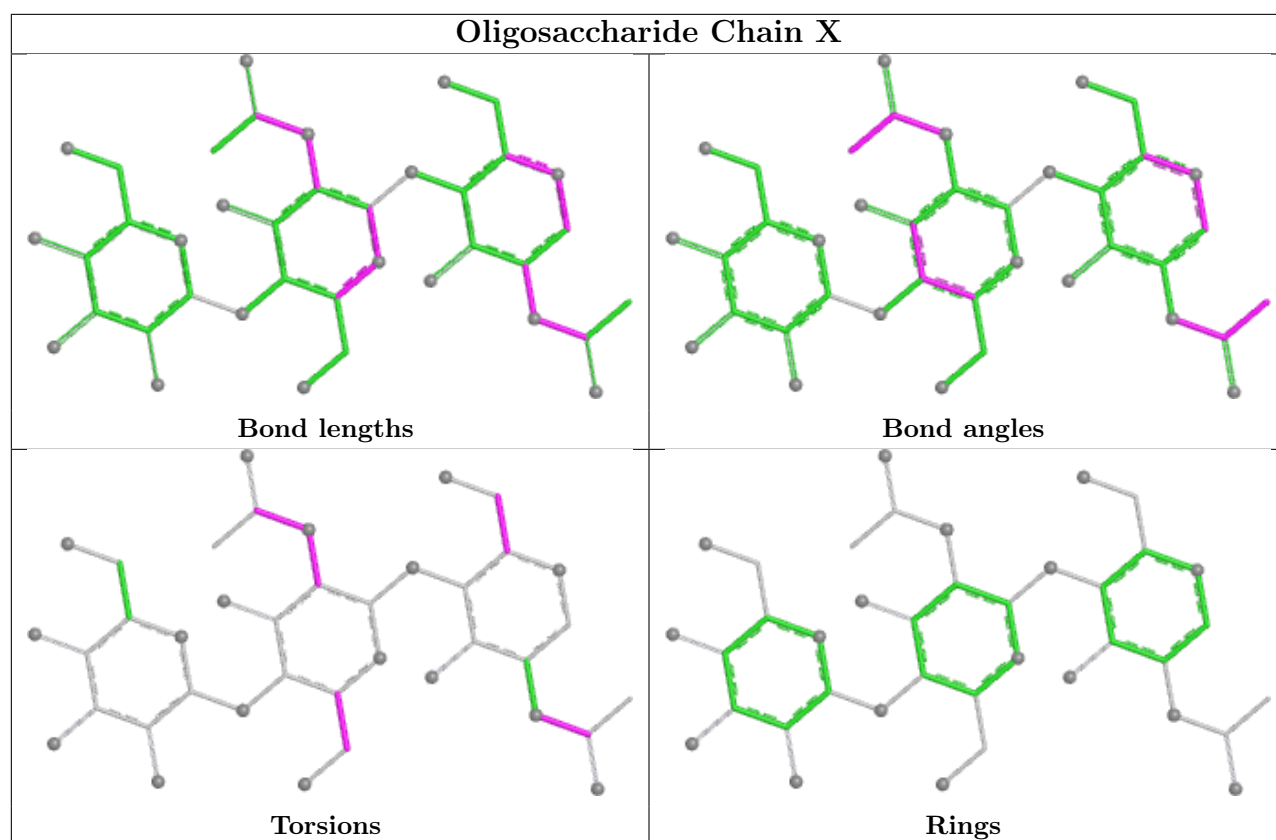


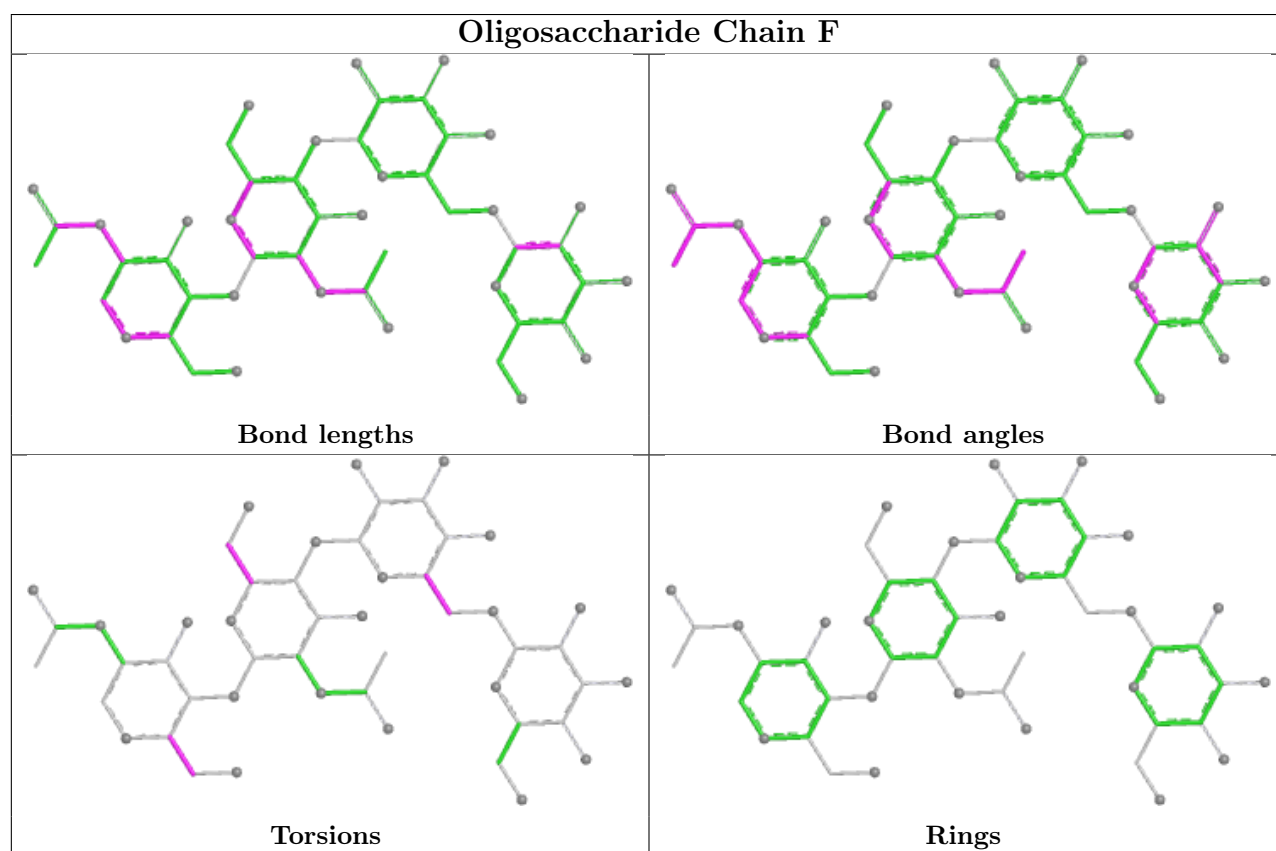
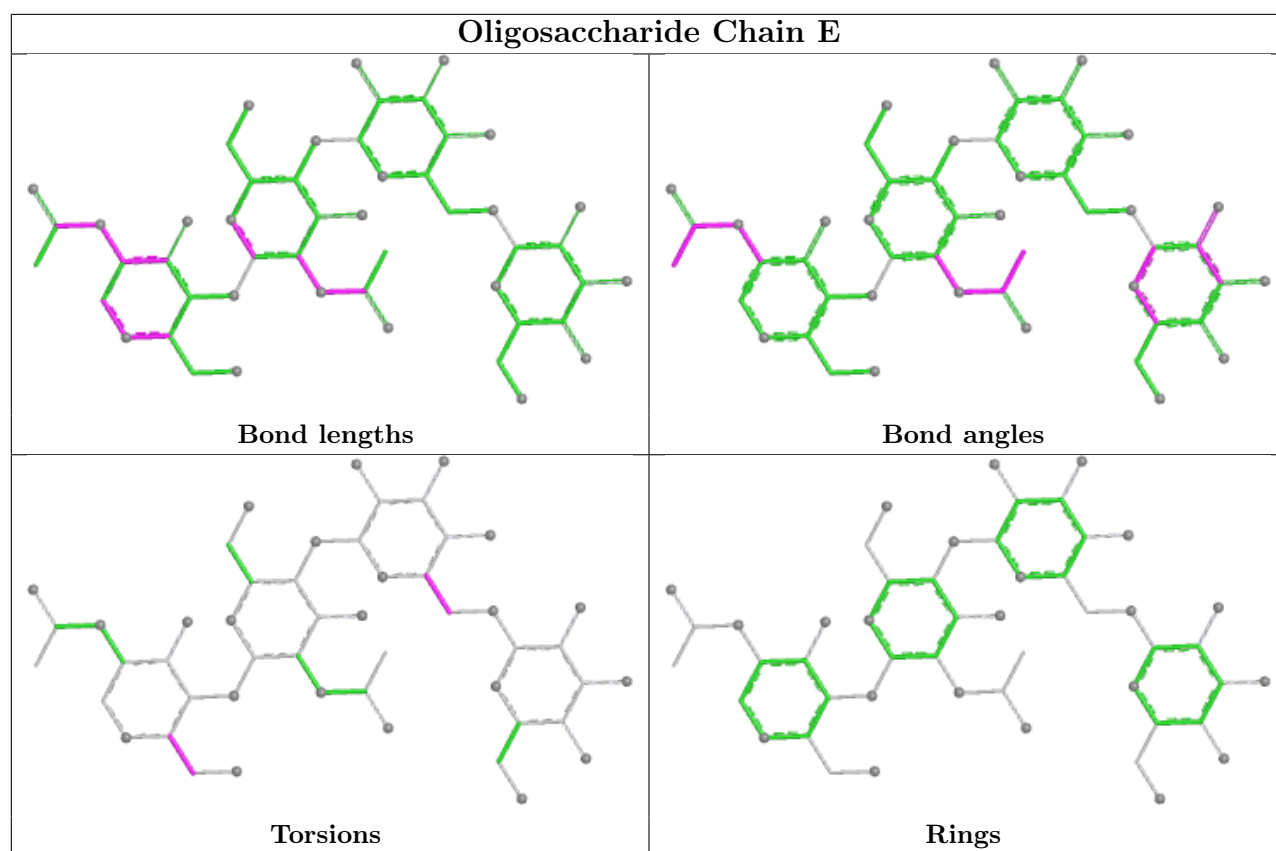


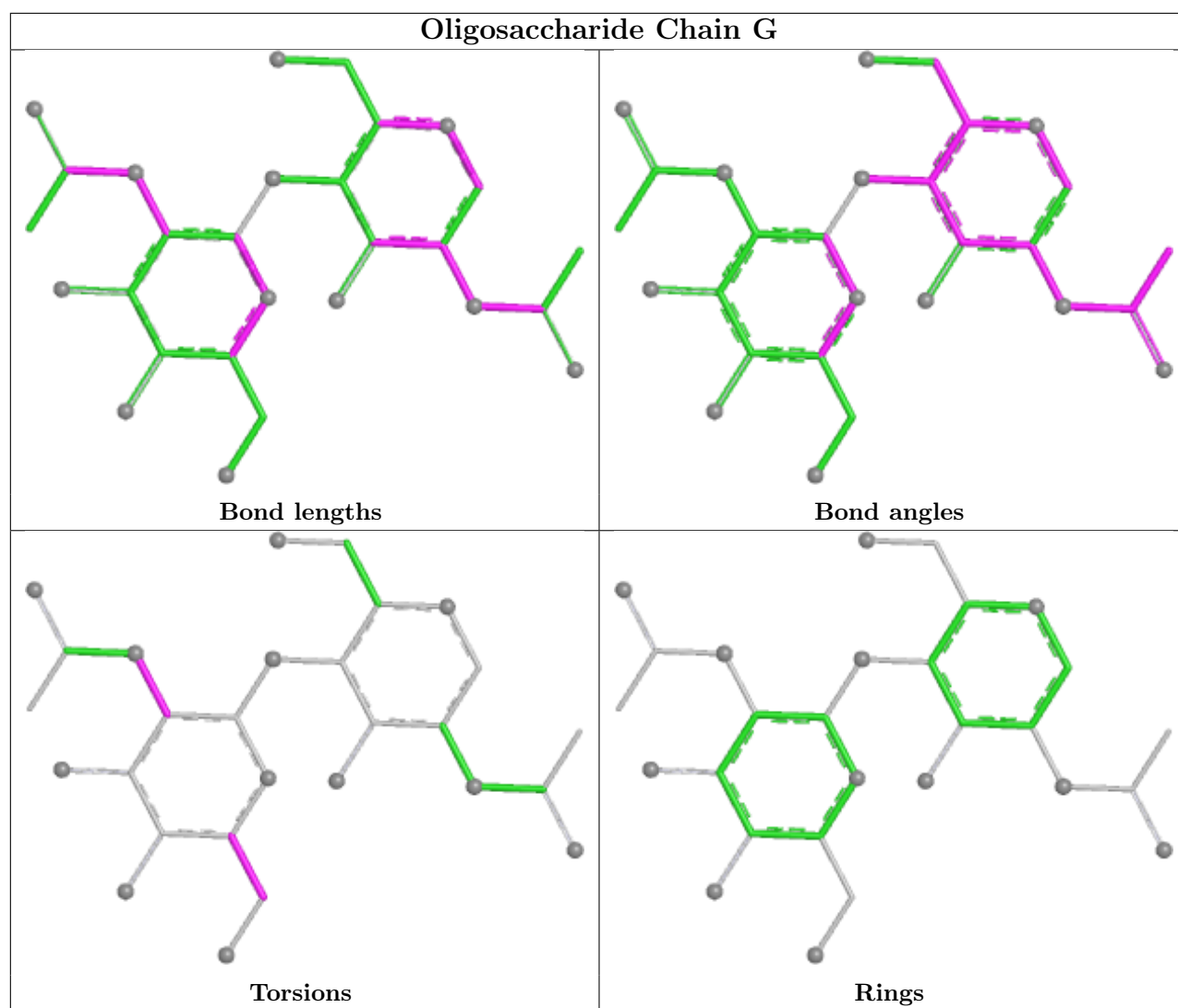


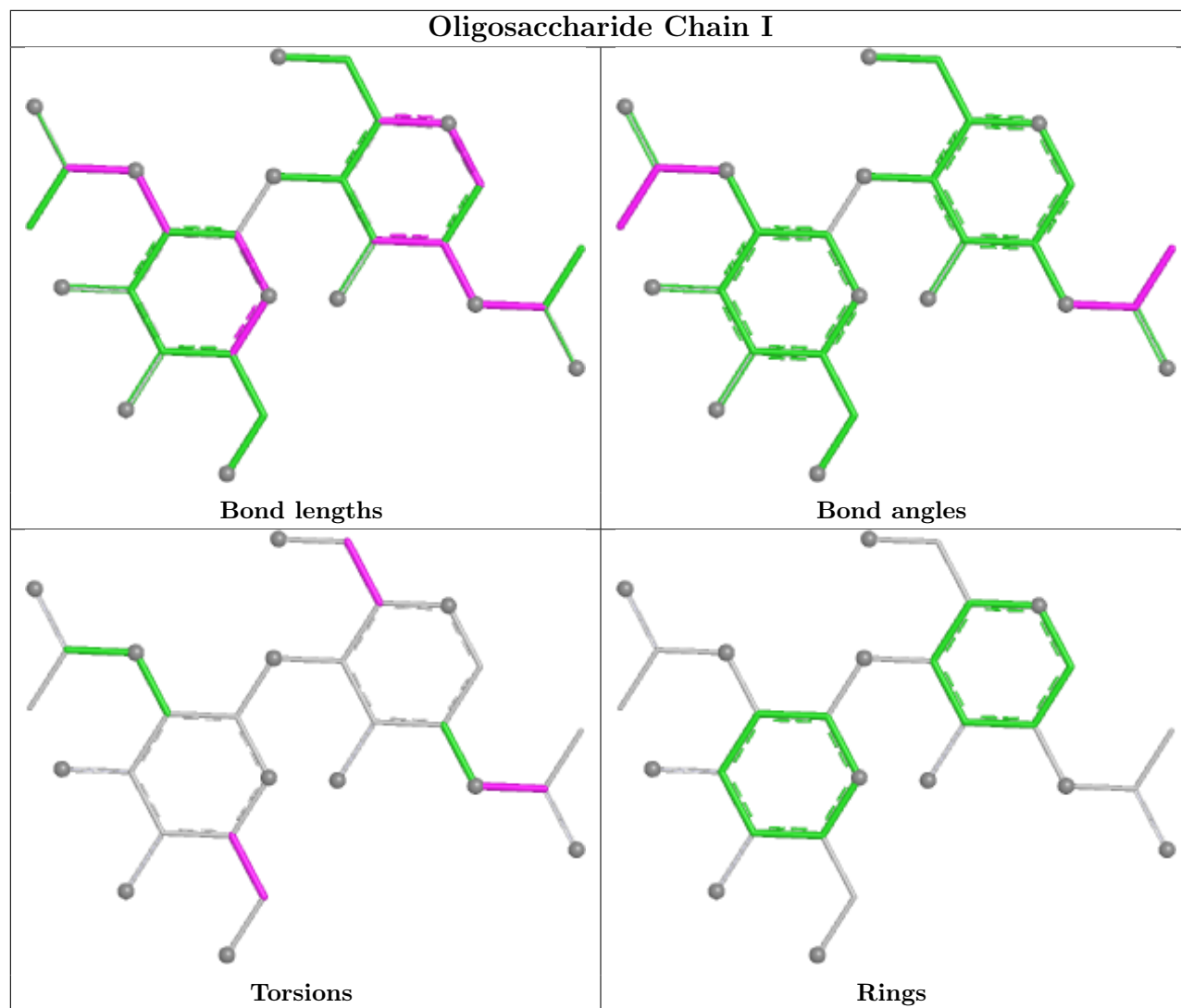


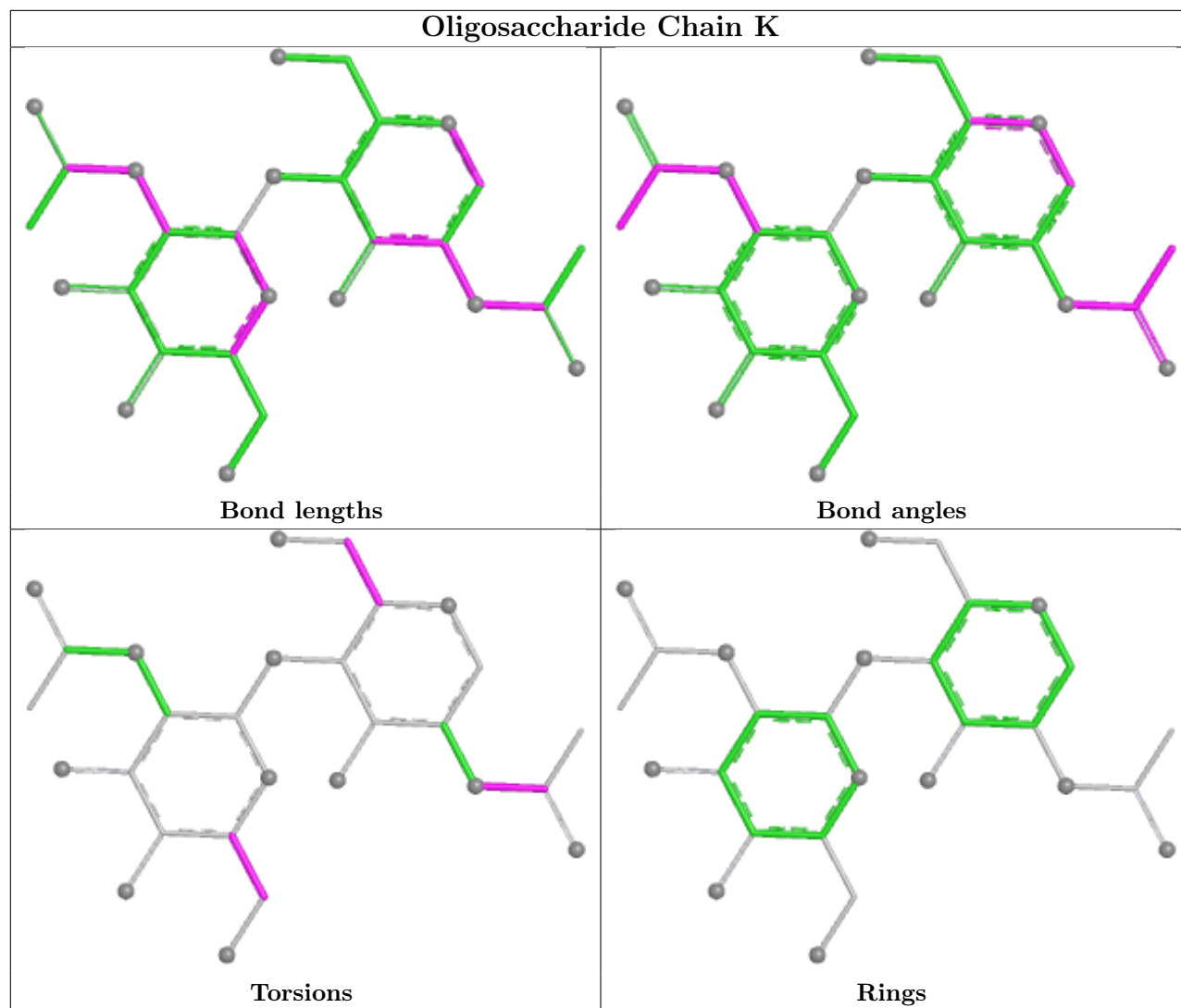


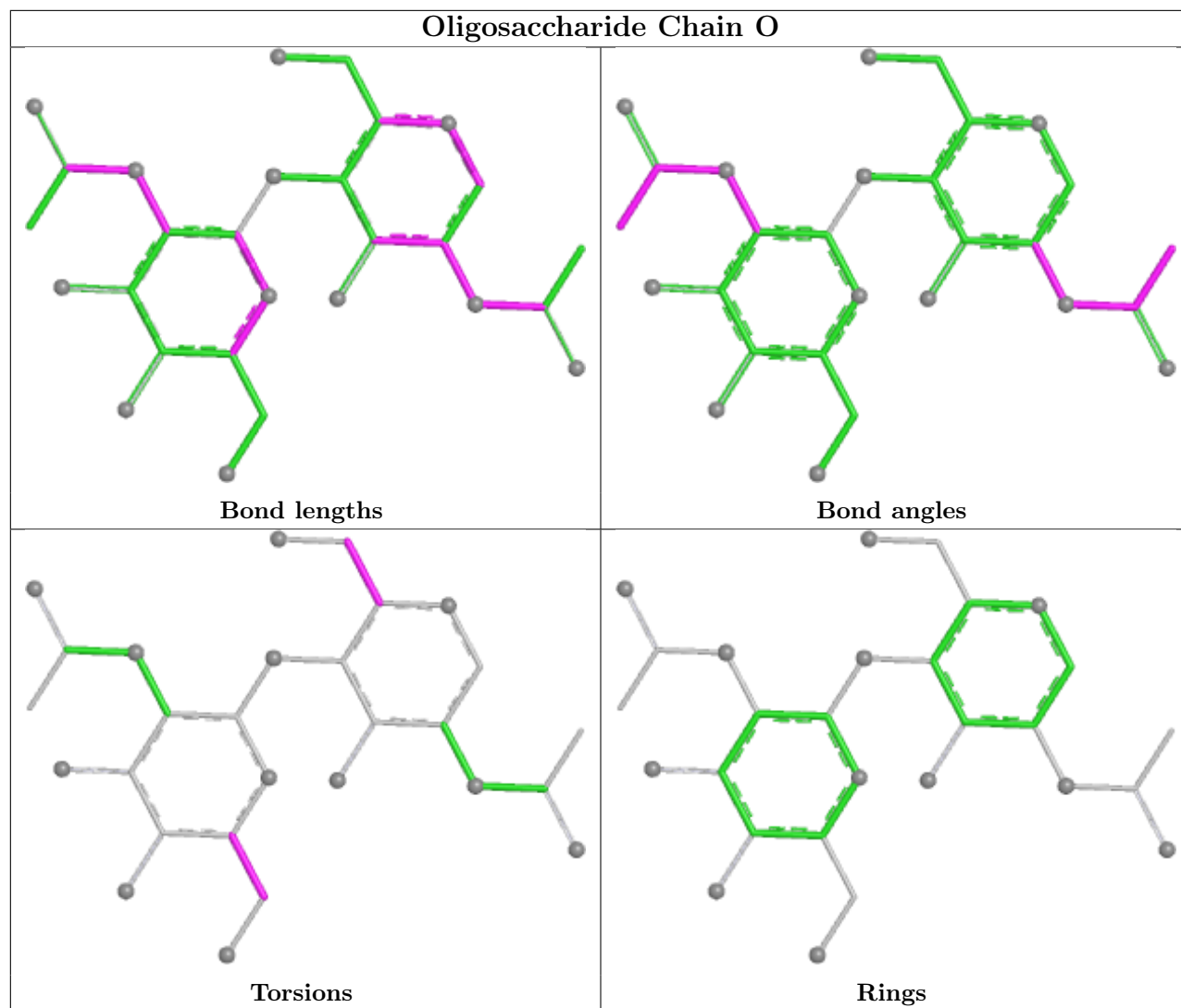


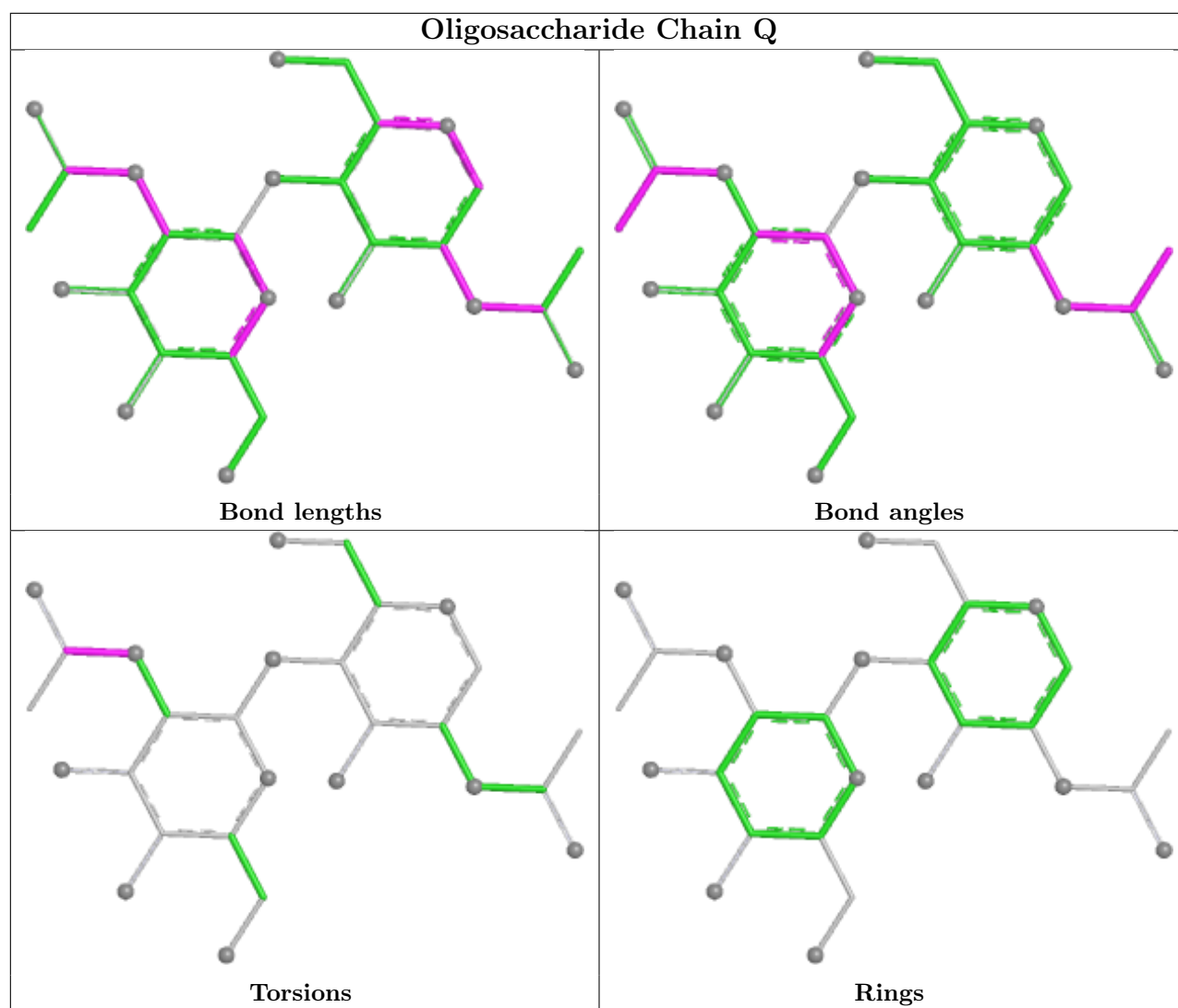




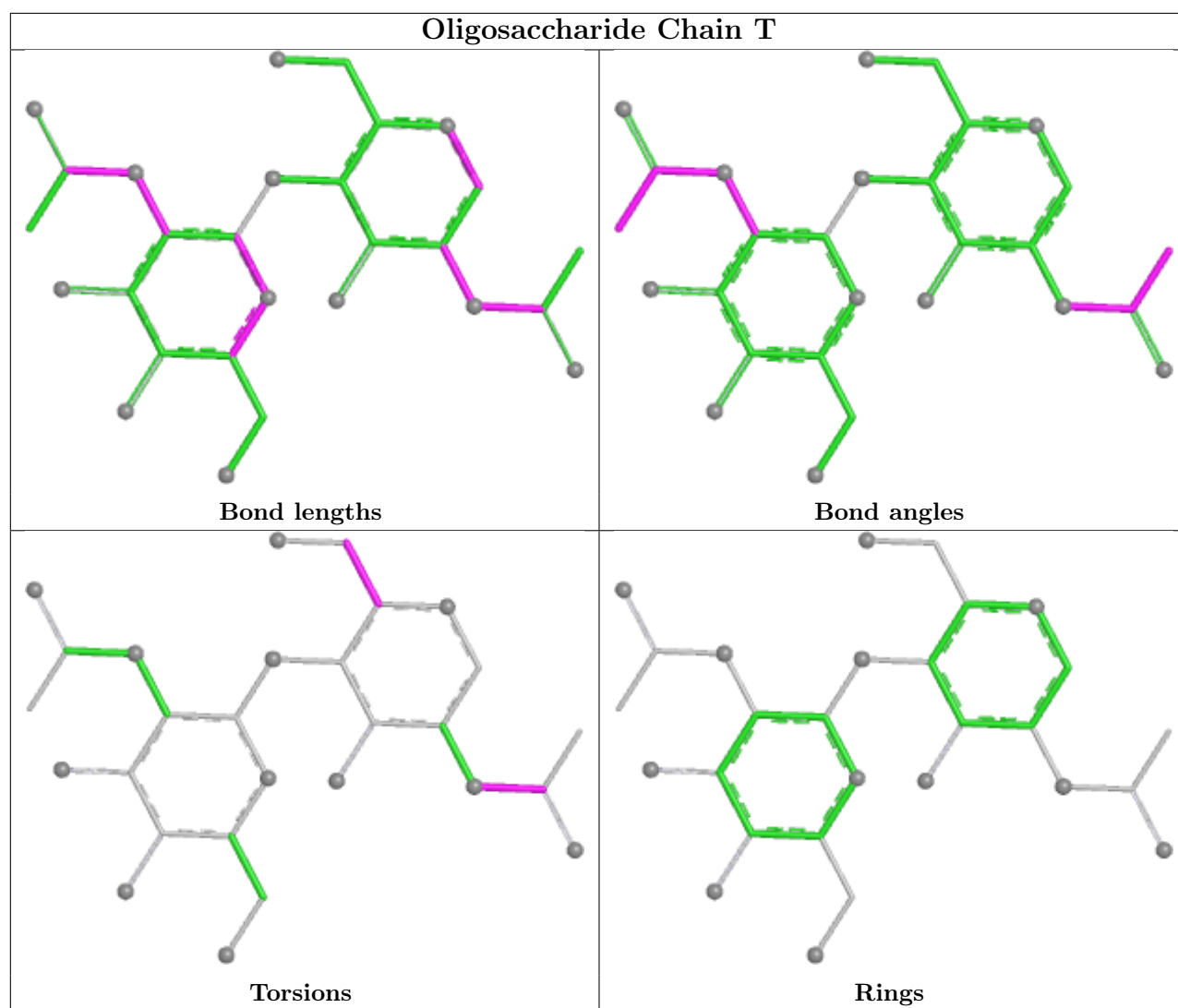


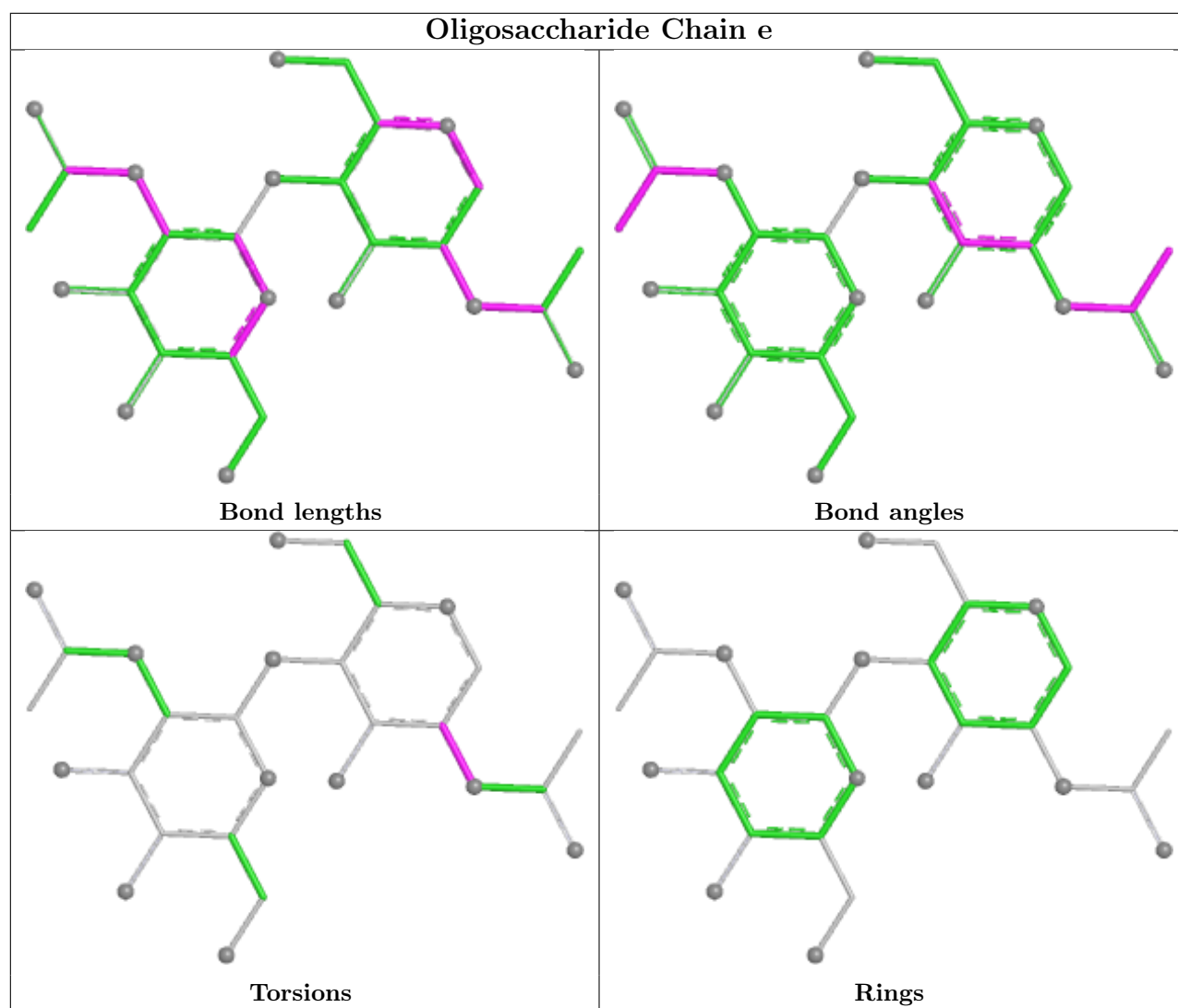


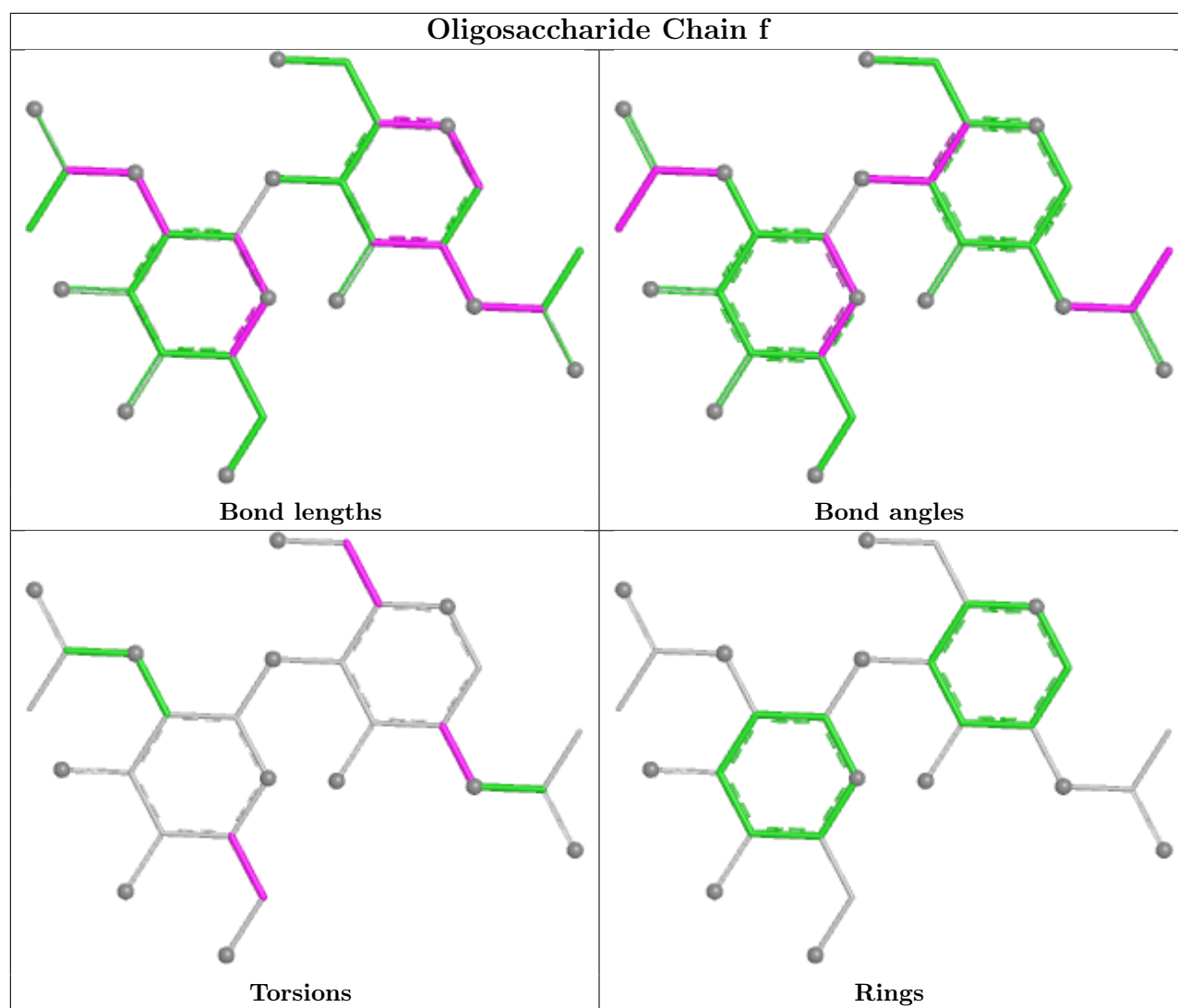


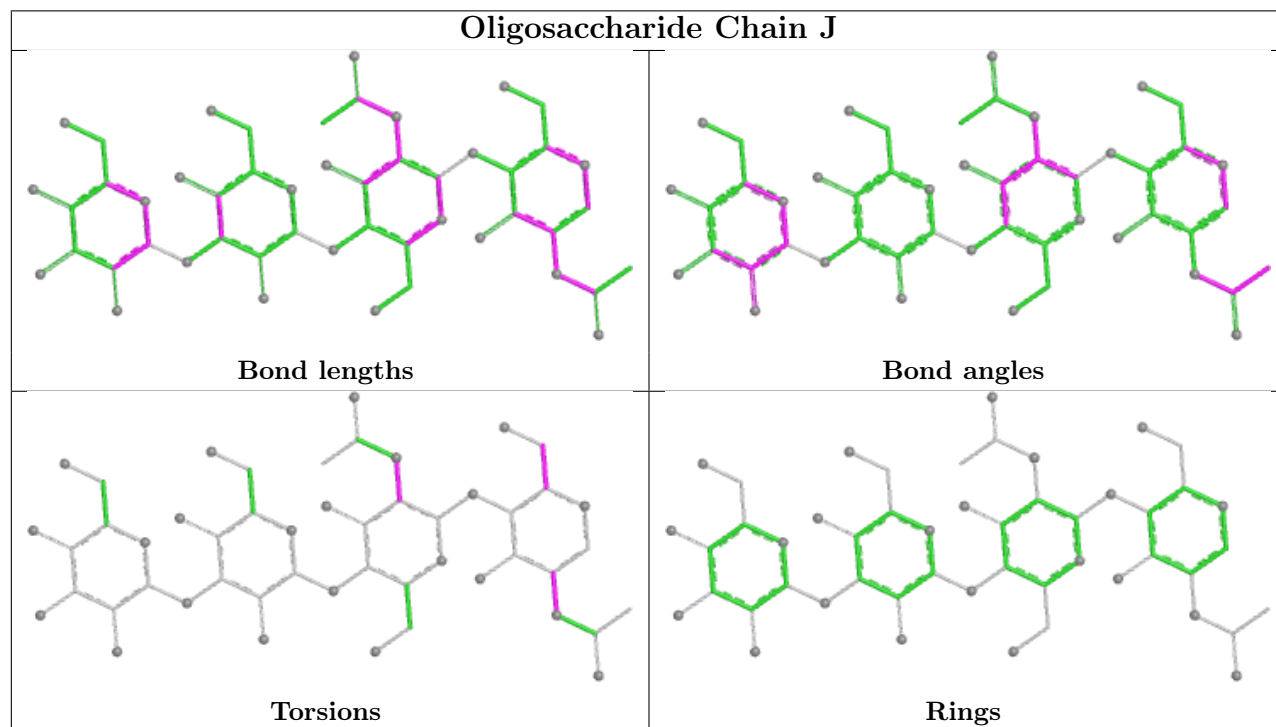
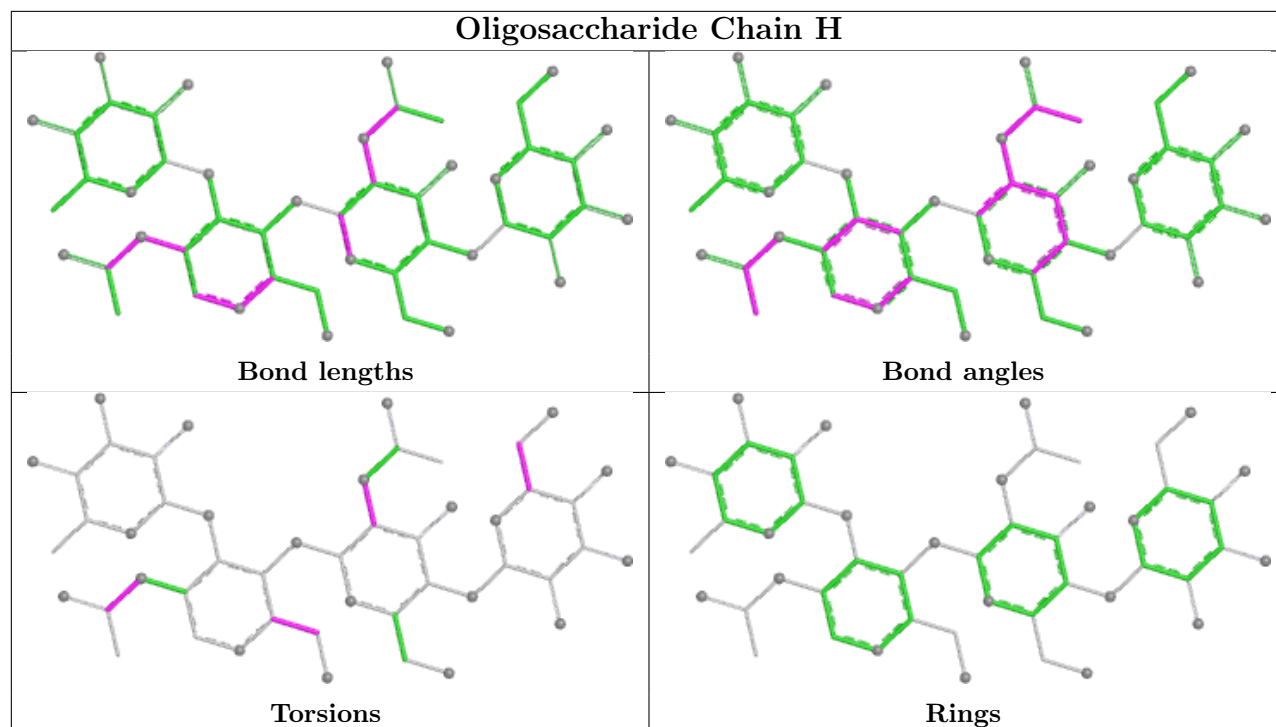


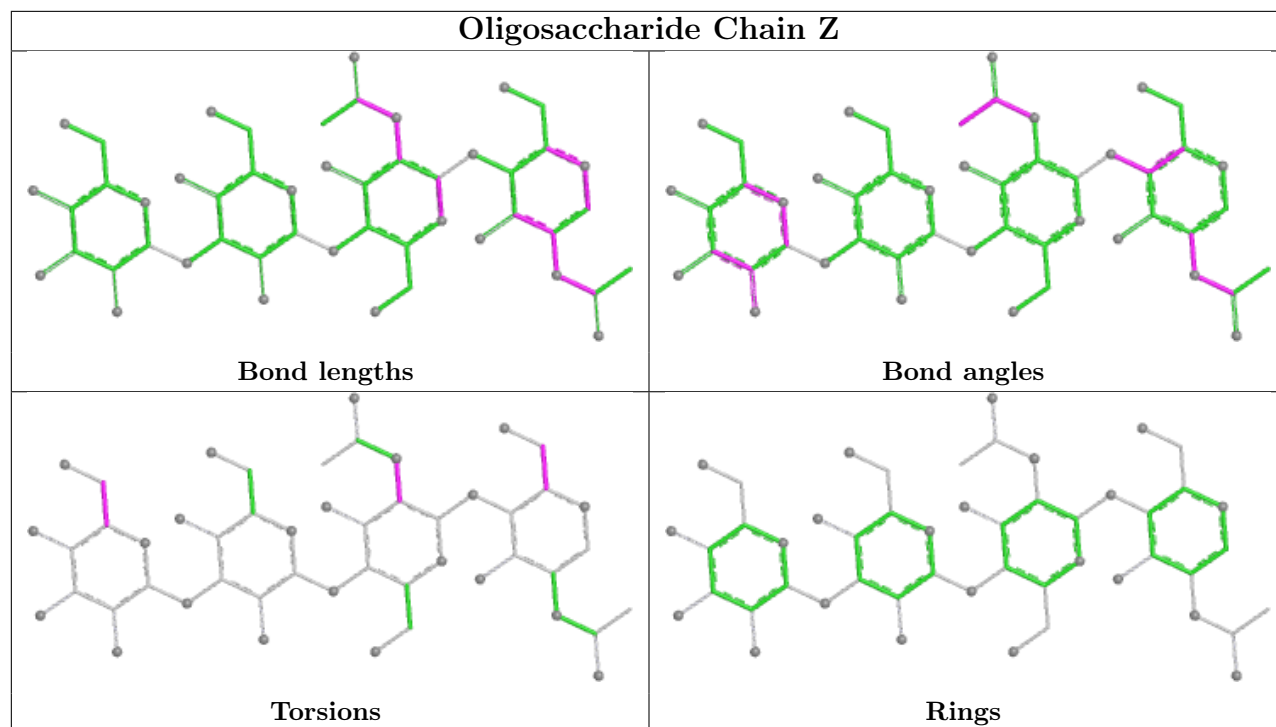


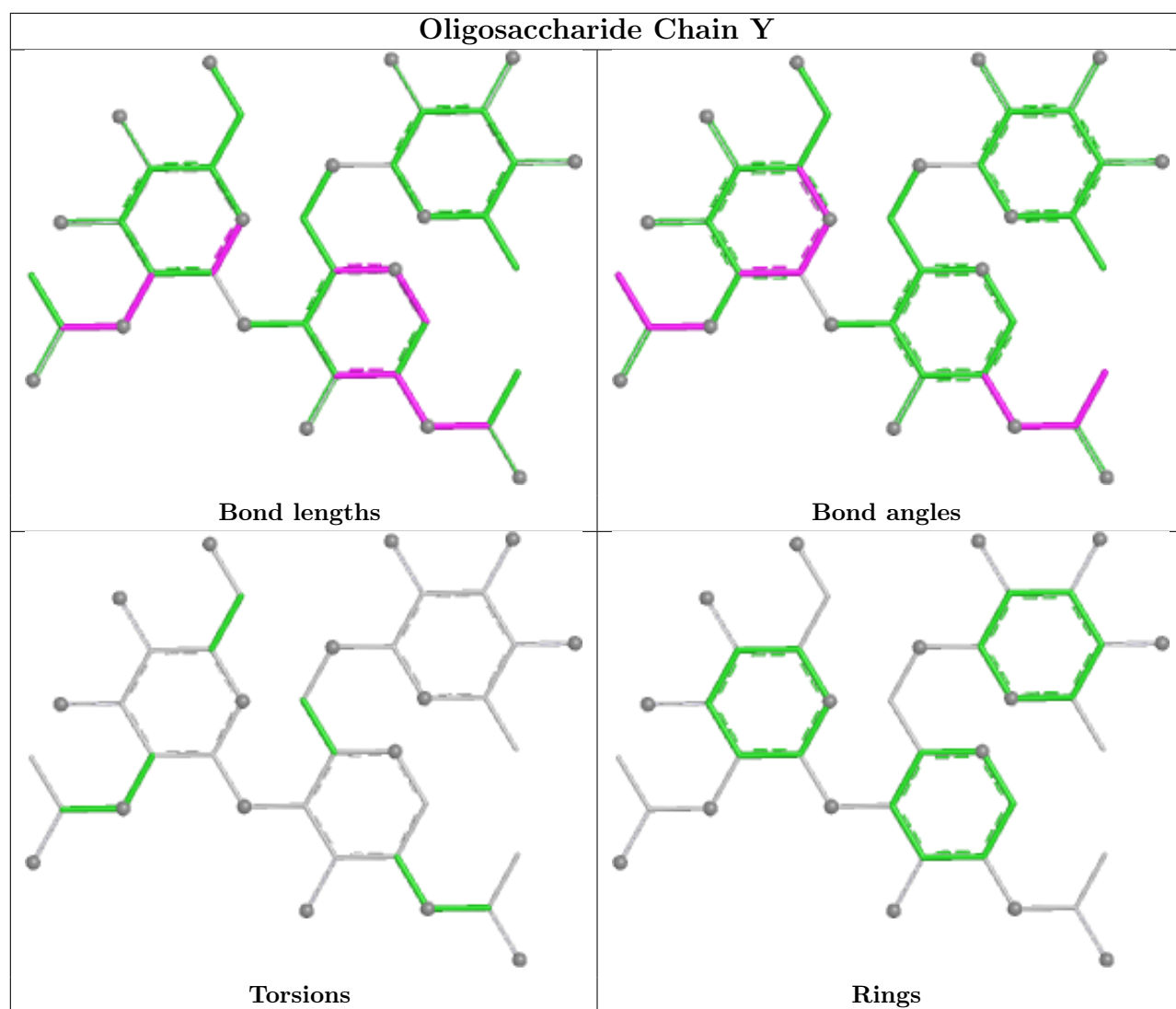


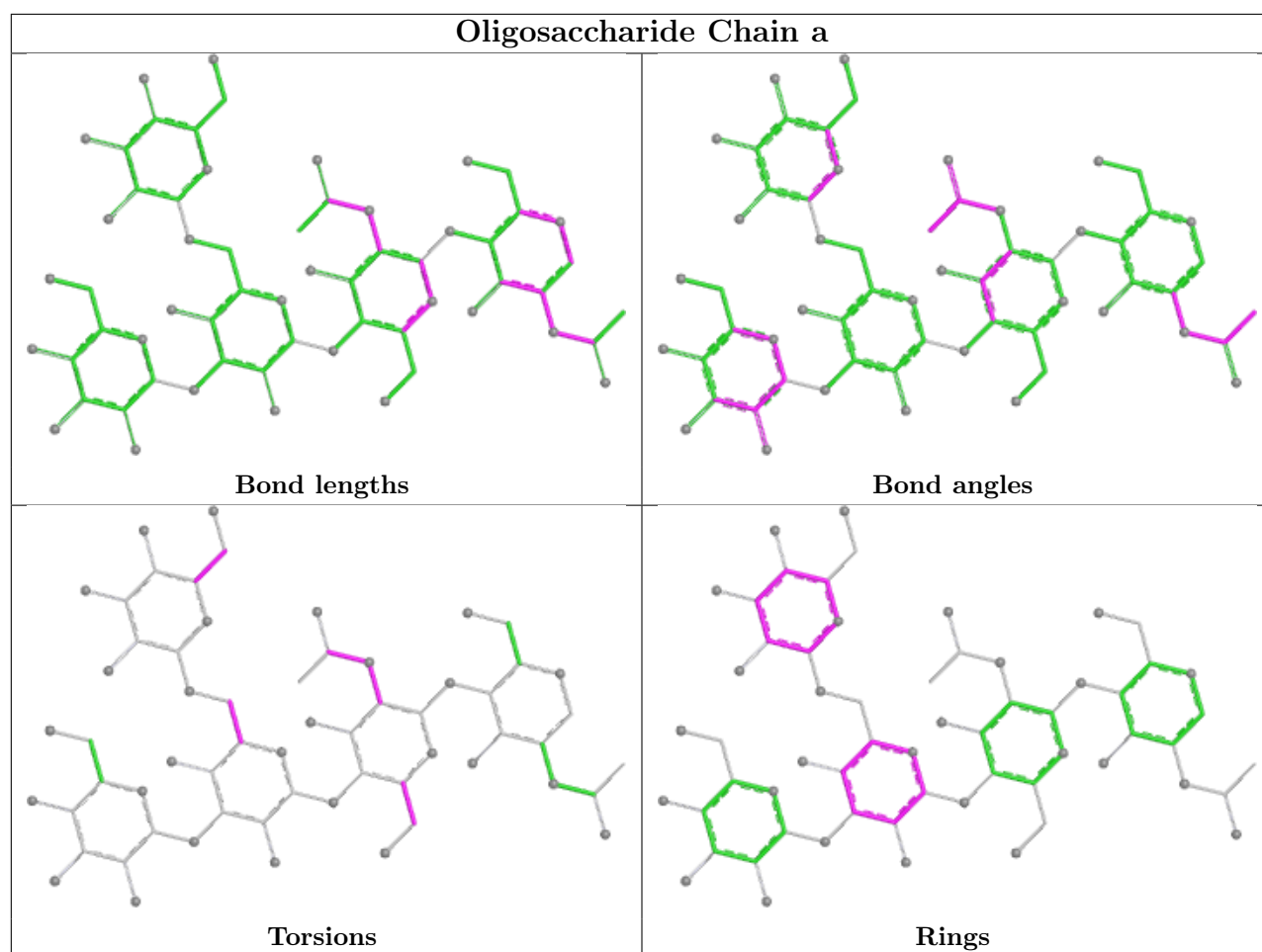


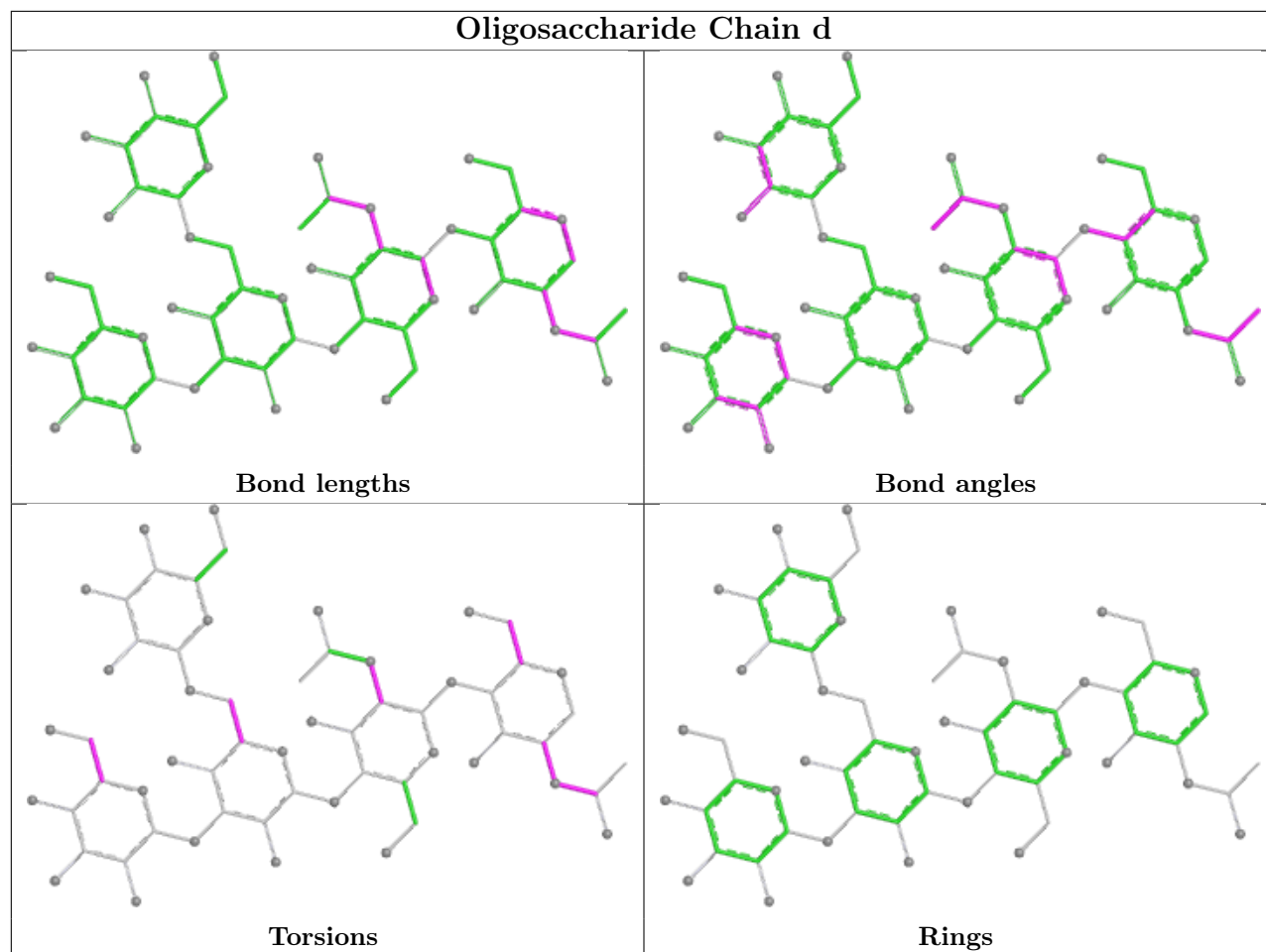




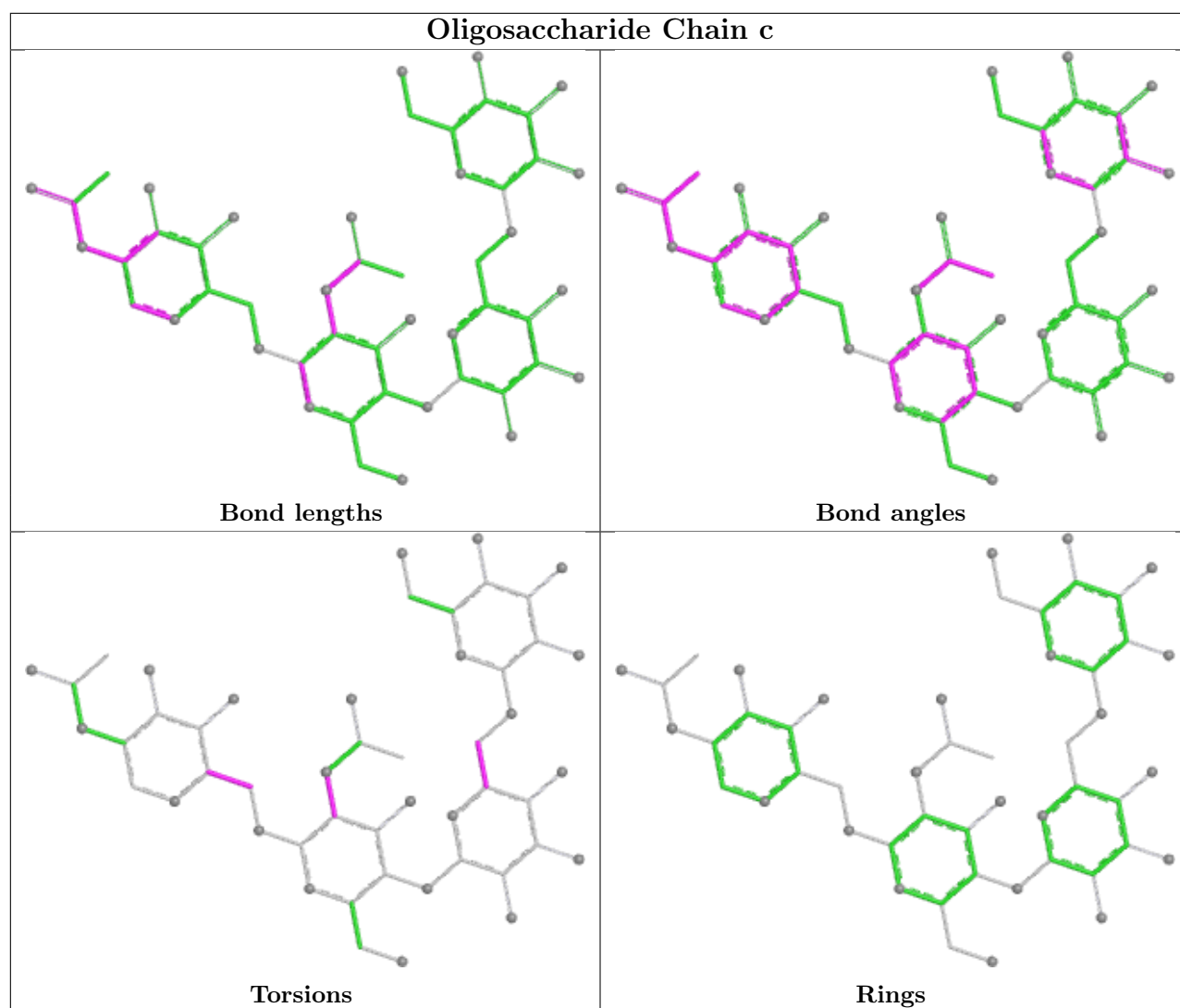












## 5.6 Ligand geometry [i](#)

18 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	$\# Z  > 2$
10	NAG	A	1302	1	14,14,15	2.02	4 (28%)	17,19,21	1.19	2 (11%)
10	NAG	C	1401	1	14,14,15	2.03	3 (21%)	17,19,21	1.50	2 (11%)
10	NAG	B	1404	1	14,14,15	1.99	4 (28%)	17,19,21	1.17	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
10	NAG	C	1402	1	14,14,15	2.02	4 (28%)	17,19,21	1.21	1 (5%)
10	NAG	C	1403	1	14,14,15	1.93	3 (21%)	17,19,21	1.91	2 (11%)
10	NAG	C	1404	1	14,14,15	2.01	4 (28%)	17,19,21	1.60	5 (29%)
10	NAG	A	1303	1	14,14,15	1.99	4 (28%)	17,19,21	1.04	2 (11%)
10	NAG	C	1405	1	14,14,15	2.02	4 (28%)	17,19,21	0.99	1 (5%)
10	NAG	B	1402	1	14,14,15	2.06	4 (28%)	17,19,21	1.26	2 (11%)
10	NAG	A	1306	1	14,14,15	2.04	3 (21%)	17,19,21	1.48	3 (17%)
10	NAG	A	1304	1	14,14,15	2.16	4 (28%)	17,19,21	2.13	3 (17%)
10	NAG	B	1401	1	14,14,15	2.04	4 (28%)	17,19,21	1.20	1 (5%)
10	NAG	B	1403	1	14,14,15	2.05	4 (28%)	17,19,21	1.37	4 (23%)
10	NAG	A	1305	1	14,14,15	2.21	4 (28%)	17,19,21	1.72	5 (29%)
10	NAG	B	1407	1	14,14,15	2.36	5 (35%)	17,19,21	1.46	2 (11%)
10	NAG	B	1405	1	14,14,15	1.96	4 (28%)	17,19,21	1.18	2 (11%)
10	NAG	A	1301	1	14,14,15	2.01	4 (28%)	17,19,21	1.08	2 (11%)
10	NAG	B	1406	1	14,14,15	2.09	4 (28%)	17,19,21	2.62	7 (41%)

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	A	1302	1	-	2/6/23/26	0/1/1/1
10	NAG	C	1401	1	-	4/6/23/26	0/1/1/1
10	NAG	B	1404	1	-	3/6/23/26	0/1/1/1
10	NAG	C	1402	1	-	4/6/23/26	0/1/1/1
10	NAG	C	1403	1	-	1/6/23/26	0/1/1/1
10	NAG	C	1404	1	-	0/6/23/26	0/1/1/1
10	NAG	A	1303	1	-	0/6/23/26	0/1/1/1
10	NAG	C	1405	1	-	4/6/23/26	0/1/1/1
10	NAG	B	1402	1	-	3/6/23/26	0/1/1/1
10	NAG	A	1306	1	-	2/6/23/26	0/1/1/1
10	NAG	A	1304	1	-	3/6/23/26	0/1/1/1
10	NAG	B	1401	1	-	2/6/23/26	0/1/1/1
10	NAG	B	1403	1	-	3/6/23/26	0/1/1/1
10	NAG	A	1305	1	-	4/6/23/26	0/1/1/1
10	NAG	B	1407	1	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	NAG	B	1405	1	-	2/6/23/26	0/1/1/1
10	NAG	A	1301	1	-	0/6/23/26	0/1/1/1
10	NAG	B	1406	1	-	5/6/23/26	0/1/1/1

The worst 5 of 70 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	A	1305	NAG	O5-C1	4.84	1.51	1.43
10	A	1304	NAG	O5-C1	4.80	1.51	1.43
10	B	1407	NAG	O5-C1	4.69	1.51	1.43
10	B	1403	NAG	O5-C1	4.53	1.51	1.43
10	B	1401	NAG	O5-C1	4.41	1.51	1.43

The worst 5 of 48 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	C	1403	NAG	C1-O5-C5	6.13	120.40	112.19
10	B	1406	NAG	C1-O5-C5	4.91	118.77	112.19
10	B	1406	NAG	C4-C3-C2	4.90	118.20	111.02
10	A	1305	NAG	C8-C7-N2	4.75	124.00	116.12
10	A	1304	NAG	O5-C1-C2	4.72	118.60	111.29

There are no chirality outliers.

5 of 43 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	A	1304	NAG	C1-C2-N2-C7
10	B	1407	NAG	C3-C2-N2-C7
10	C	1405	NAG	C1-C2-N2-C7
10	B	1402	NAG	O5-C5-C6-O6
10	B	1404	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

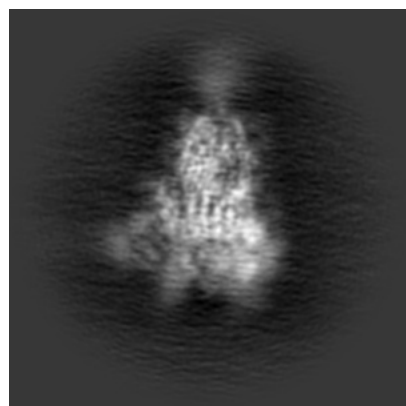
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-44344. These allow visual inspection of the internal detail of the map and identification of artifacts.

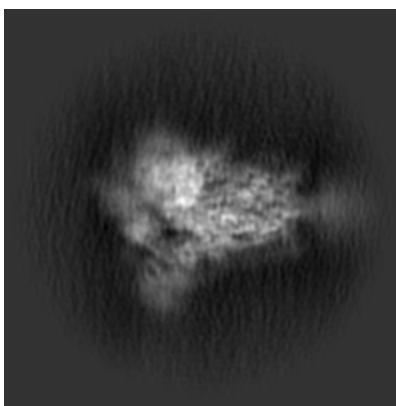
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

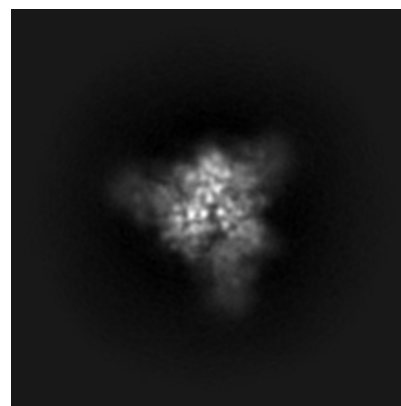
#### 6.1.1 Primary map



X

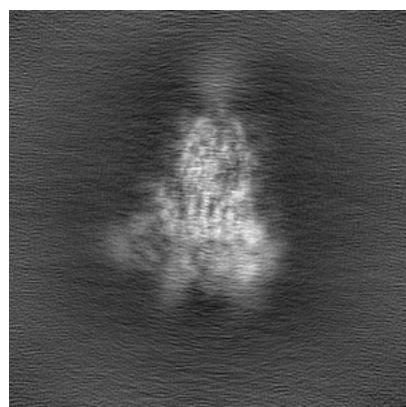


Y

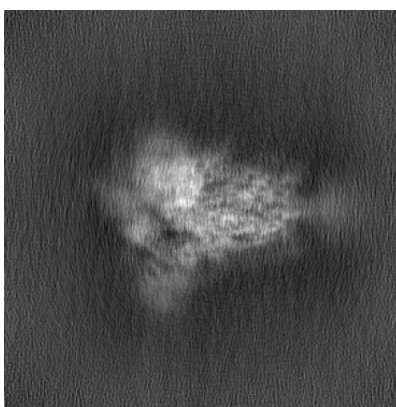


Z

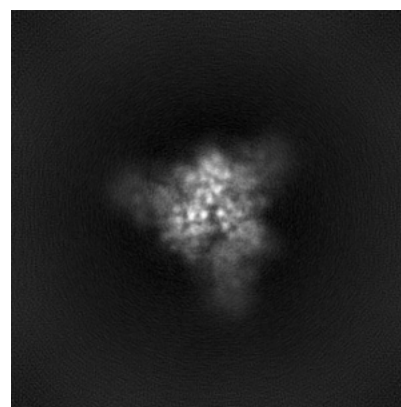
#### 6.1.2 Raw map



X



Y

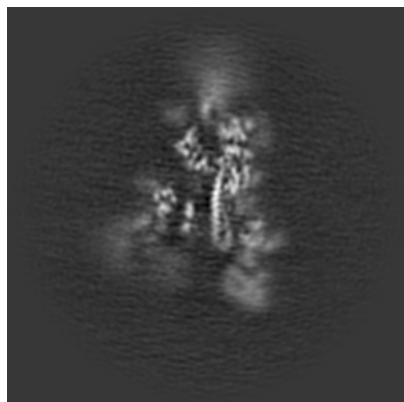


Z

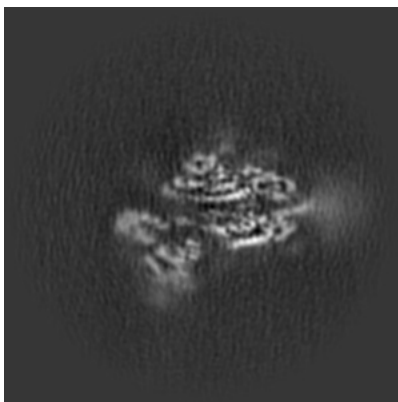
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

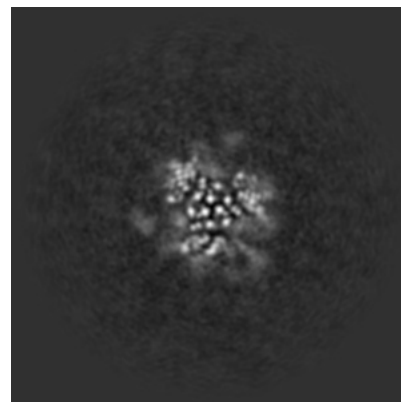
### 6.2.1 Primary map



X Index: 136

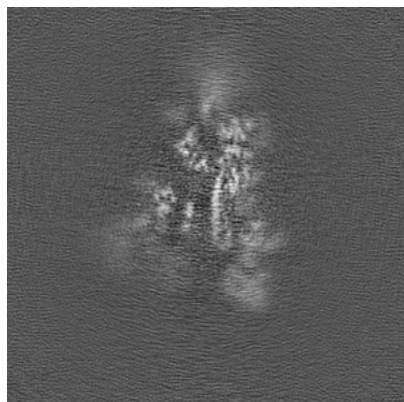


Y Index: 136

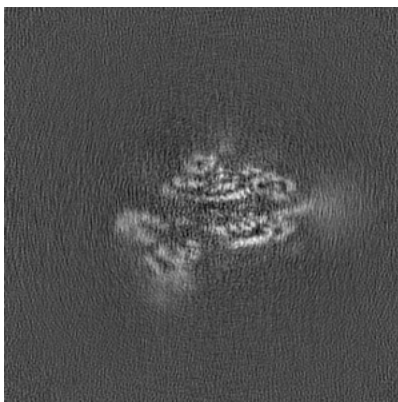


Z Index: 136

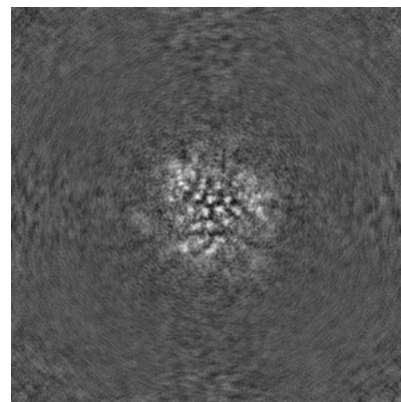
### 6.2.2 Raw map



X Index: 136



Y Index: 136

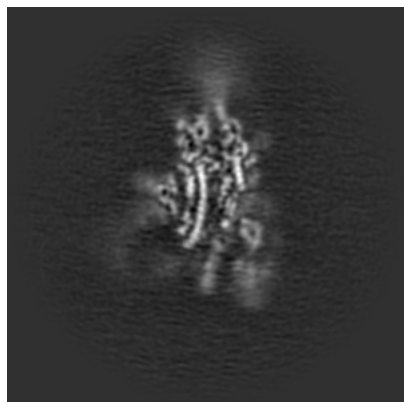


Z Index: 136

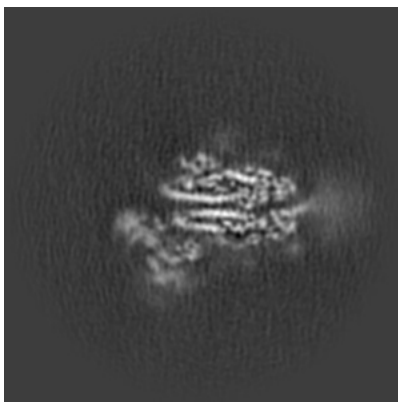
The images above show central slices of the map in three orthogonal directions.

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

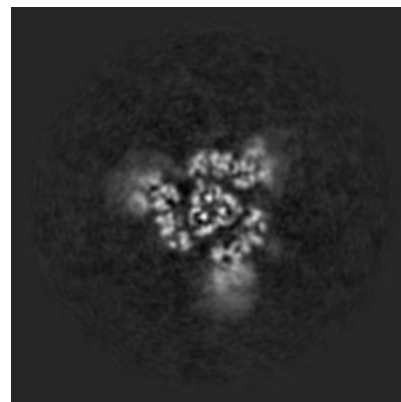
### 6.3.1 Primary map



X Index: 130

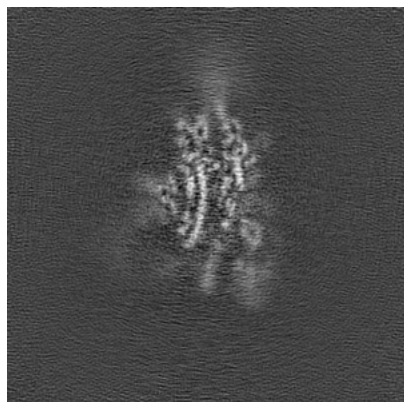


Y Index: 133

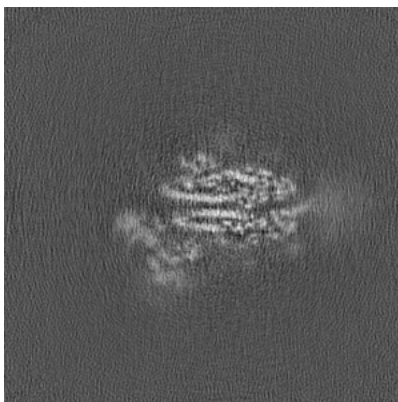


Z Index: 121

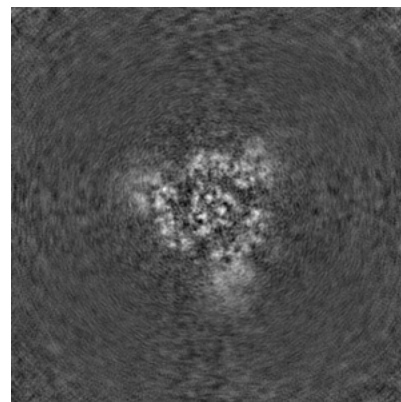
### 6.3.2 Raw map



X Index: 130



Y Index: 133



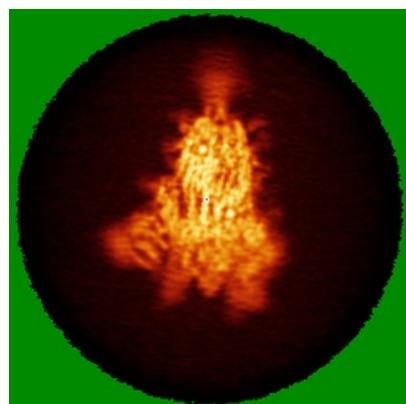
Z Index: 121

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

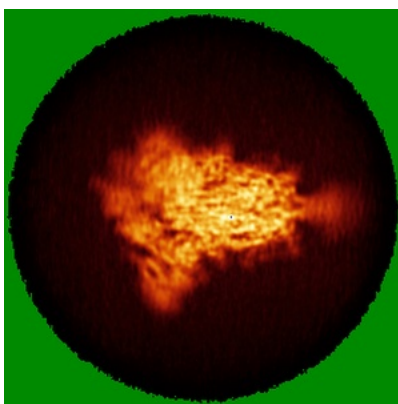


## 6.4 Orthogonal standard-deviation projections (False-color) [i](#)

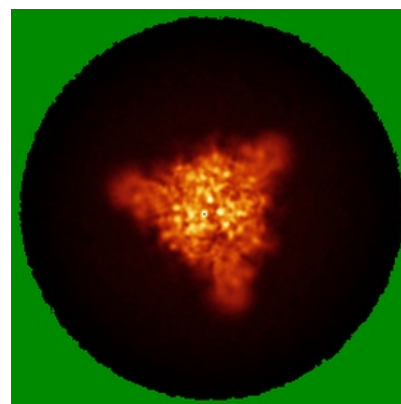
### 6.4.1 Primary map



X

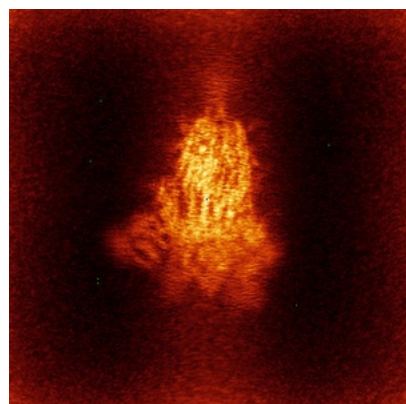


Y

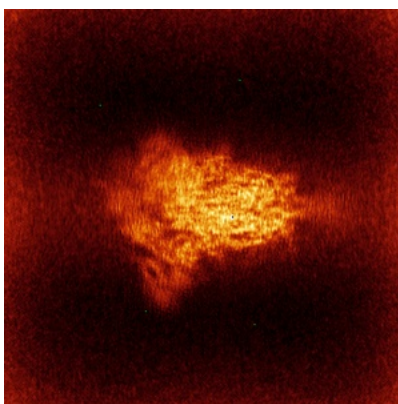


Z

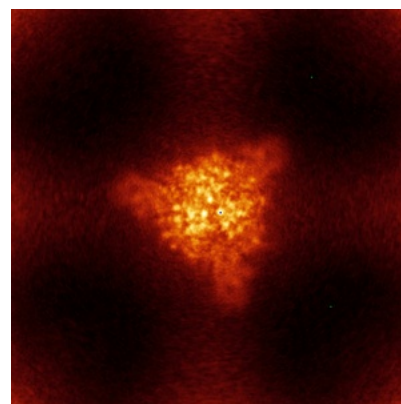
### 6.4.2 Raw map



X



Y



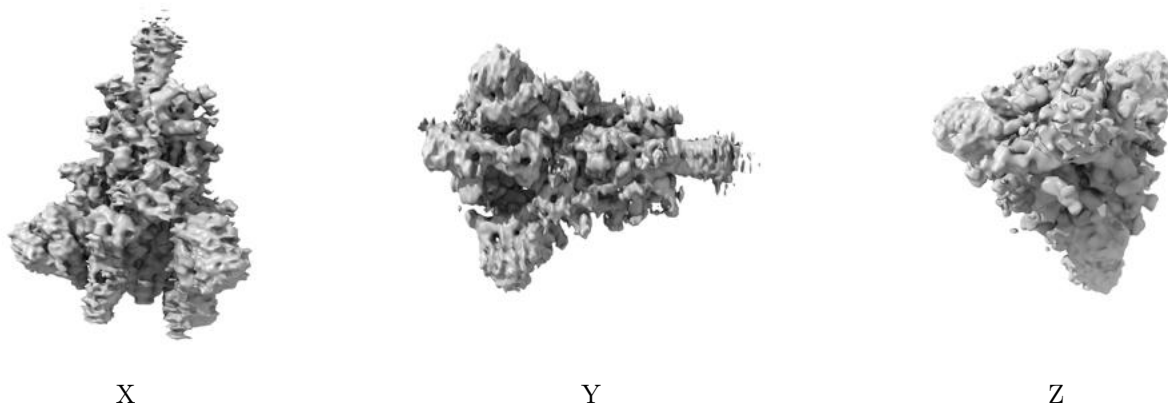
Z

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.



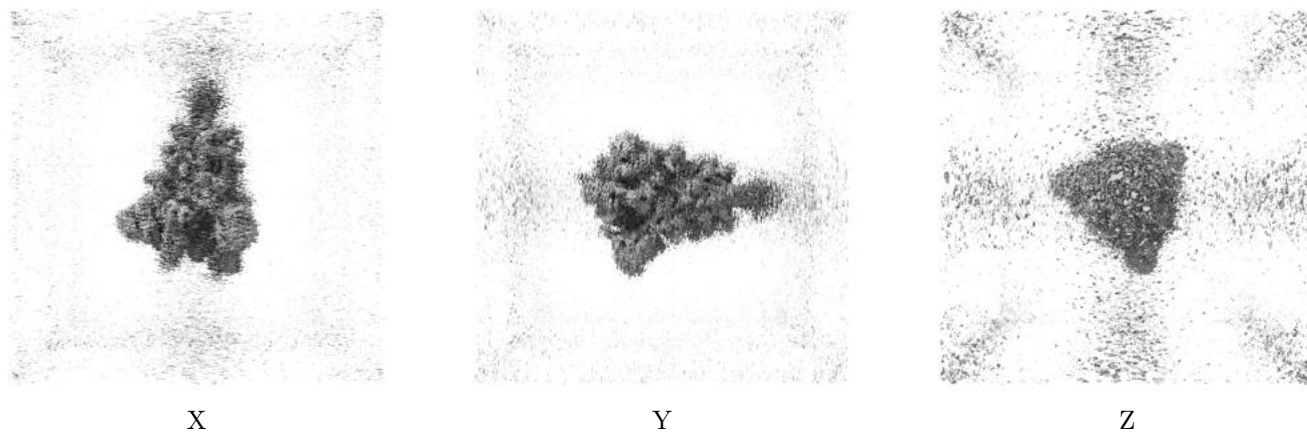
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0607. 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.

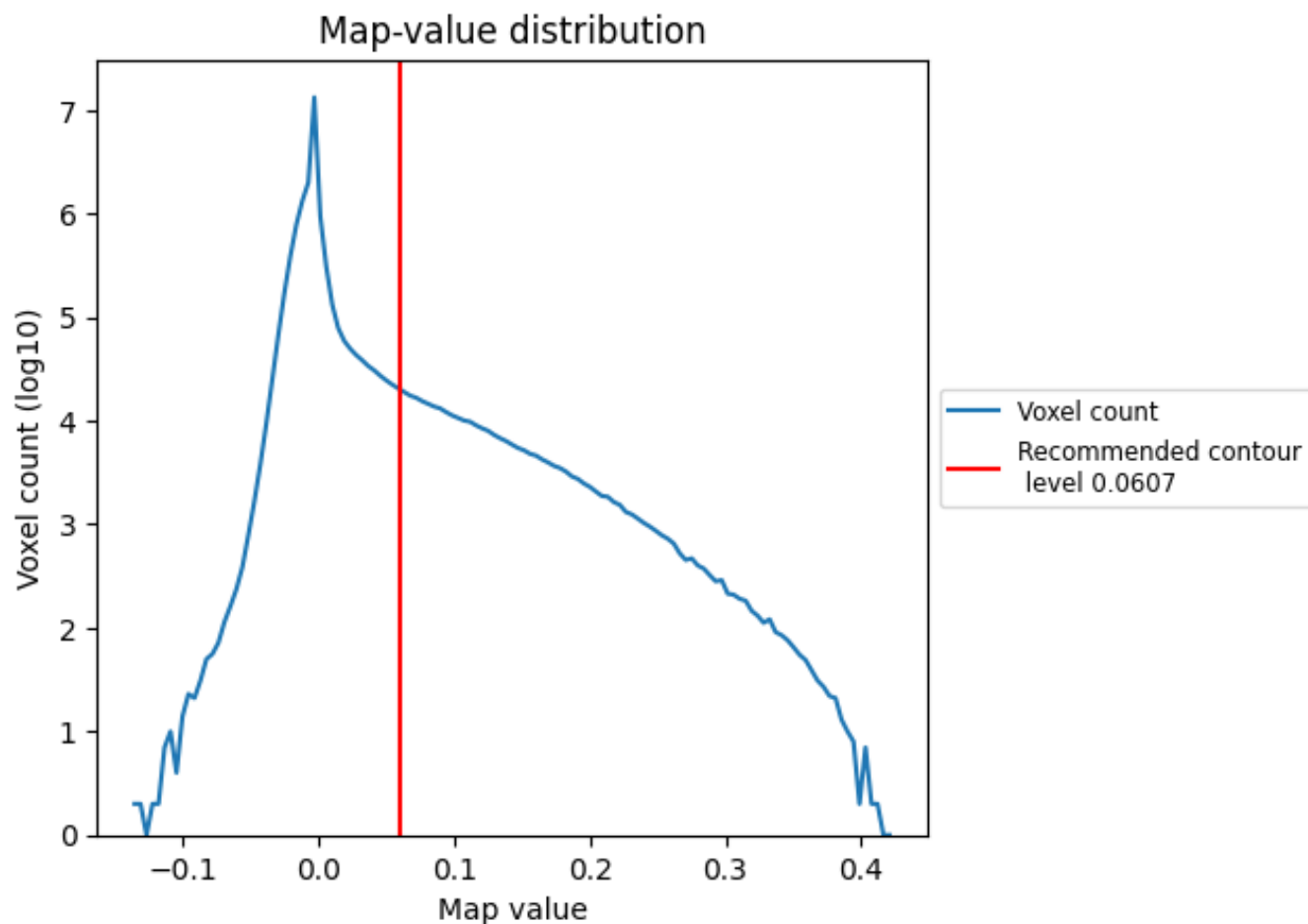
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

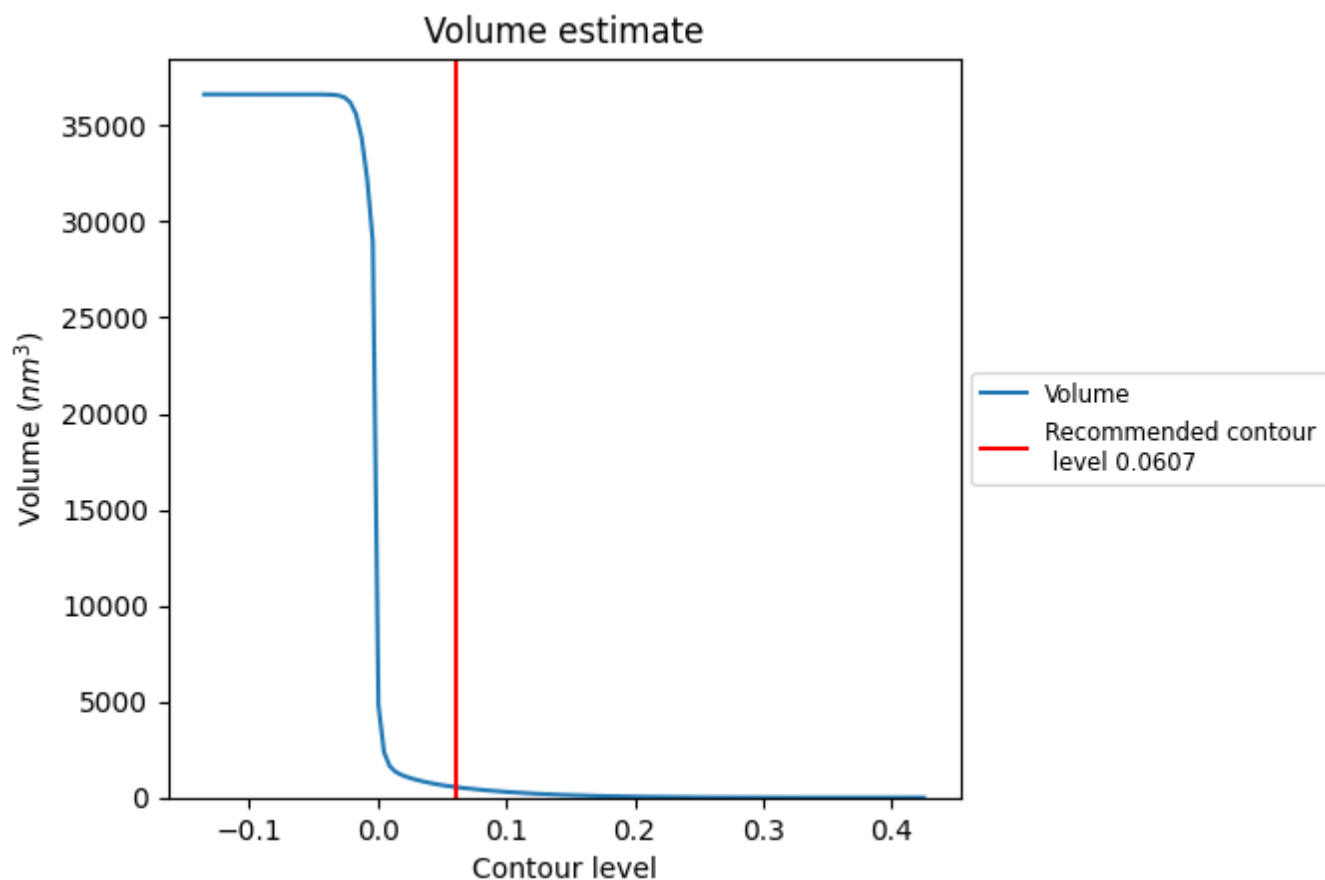
This section contains the results of statistical analysis of the map.

### 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.

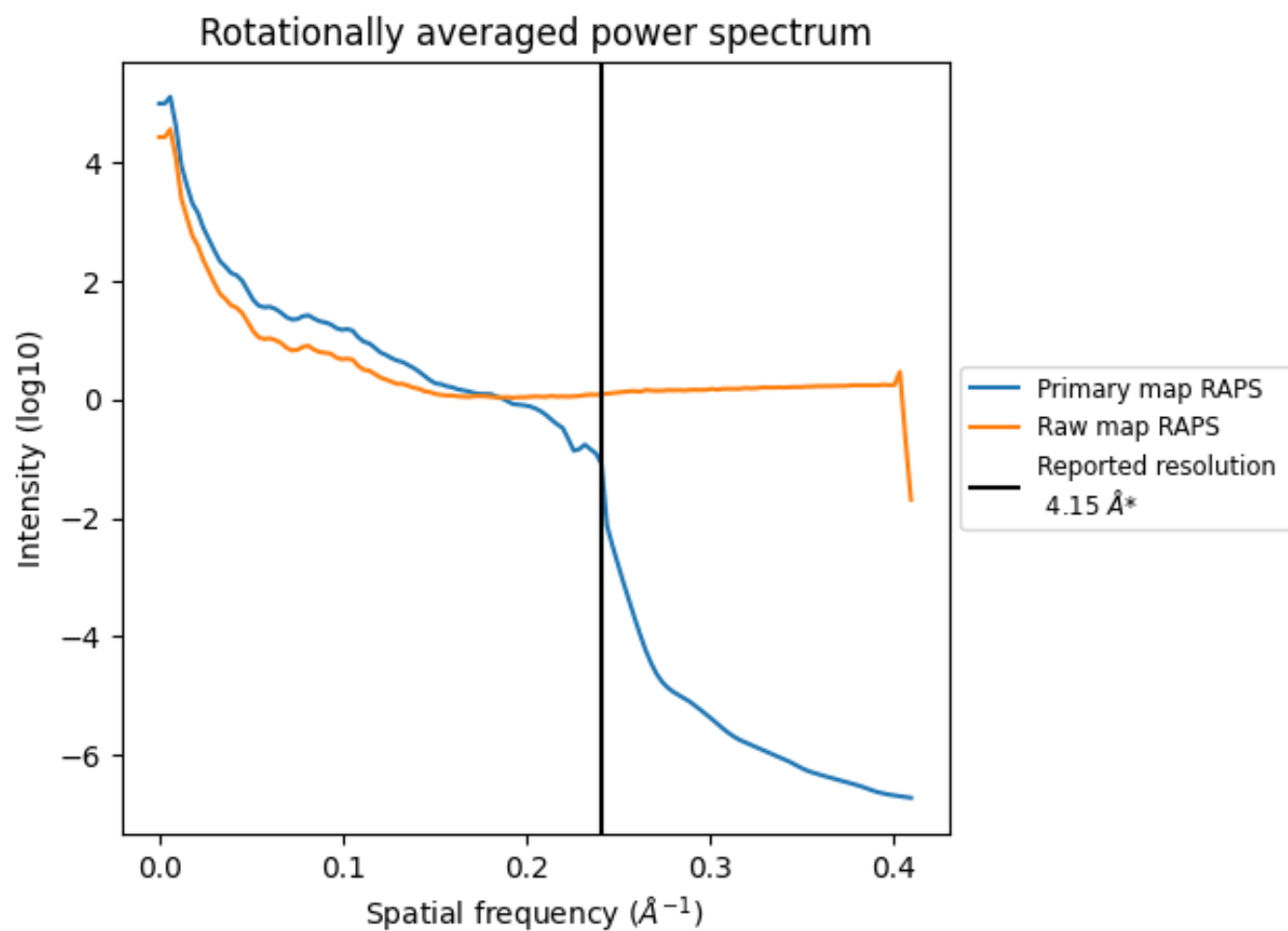
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 547 nm<sup>3</sup>; this corresponds to an approximate mass of 494 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 ⓘ

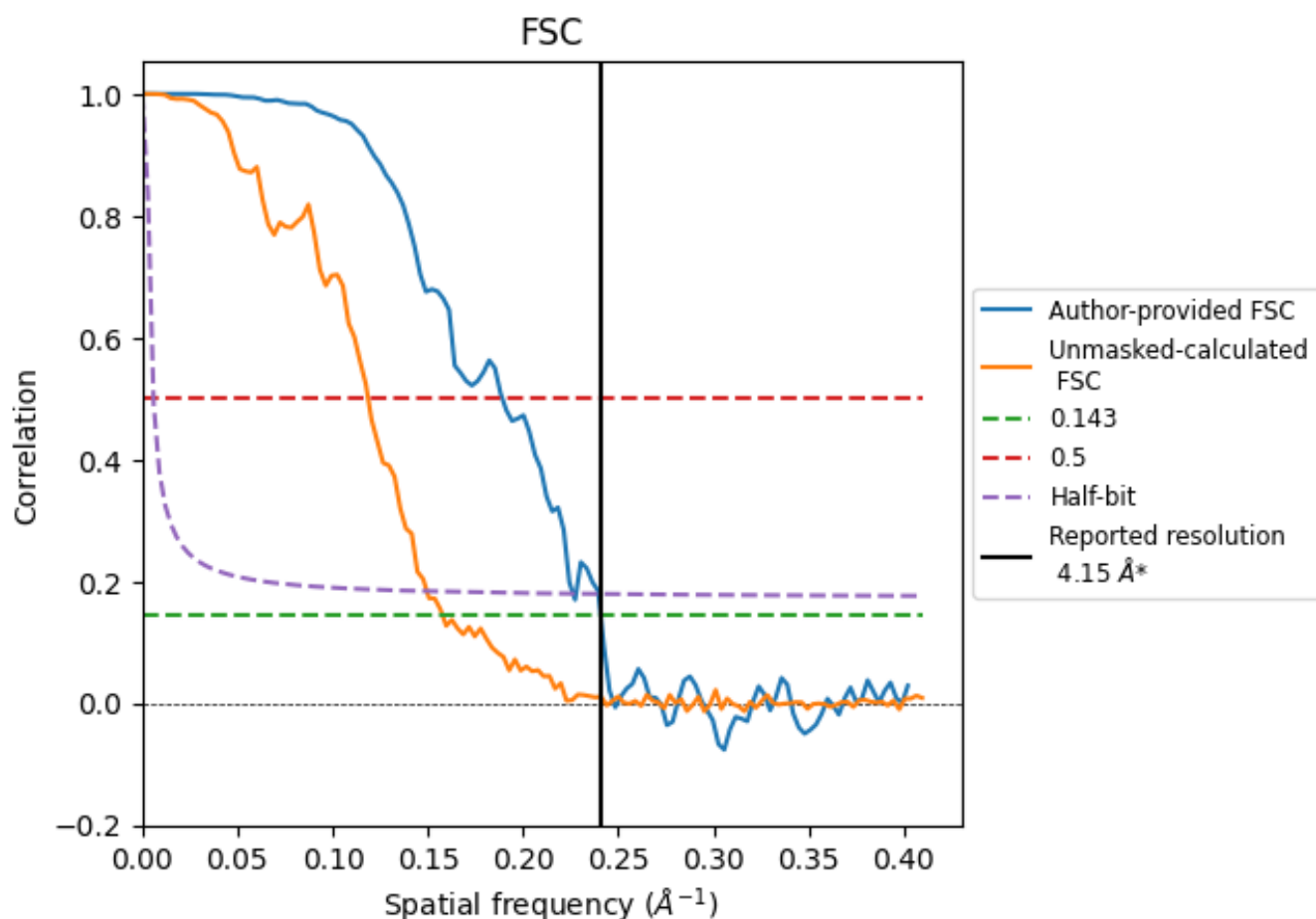


\*Reported resolution corresponds to spatial frequency of 0.241  $\text{\AA}^{-1}$

## 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 0.241  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

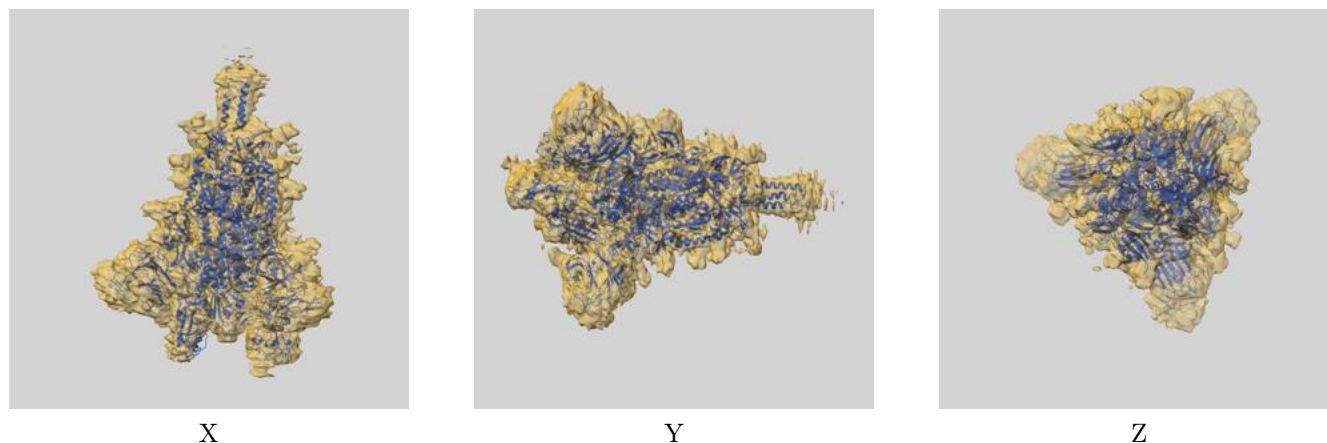
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.15	-	-
Author-provided FSC curve	4.15	5.29	4.42
Unmasked-calculated*	6.34	8.43	6.69

\*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 6.34 differs from the reported value 4.15 by more than 10 %

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

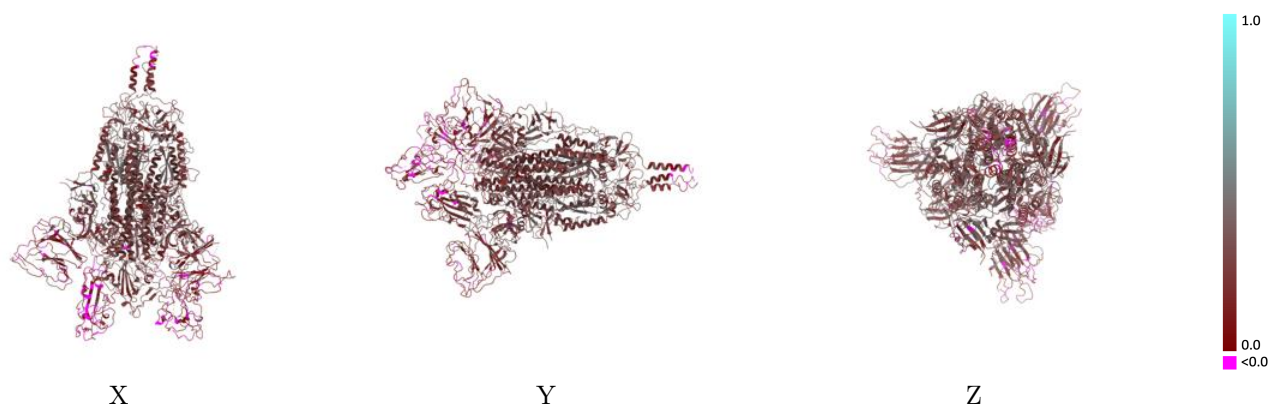
This section contains information regarding the fit between EMDB map EMD-44344 and PDB model 9B8F. Per-residue inclusion information can be found in section [3](#) on page [13](#).

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



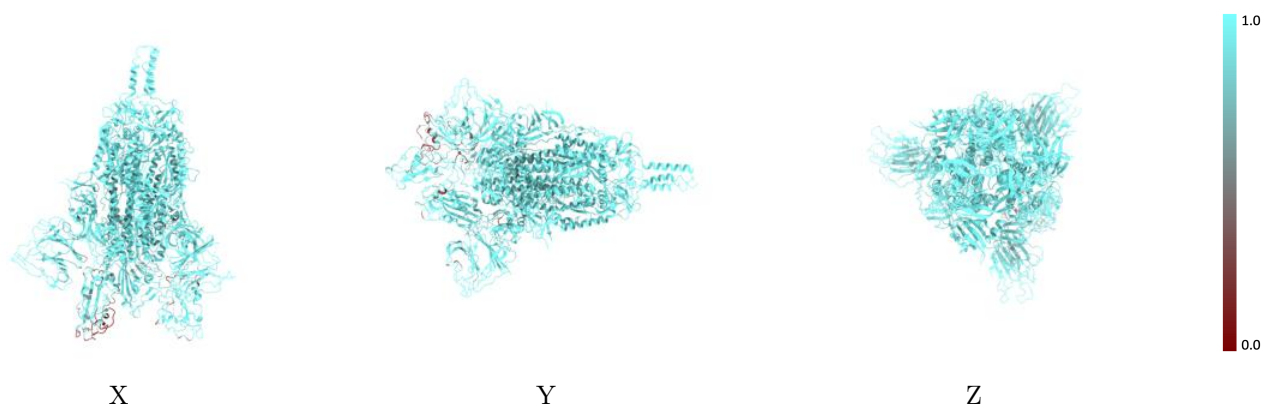
The images above show the 3D surface view of the map at the recommended contour level 0.0607 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

## 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.

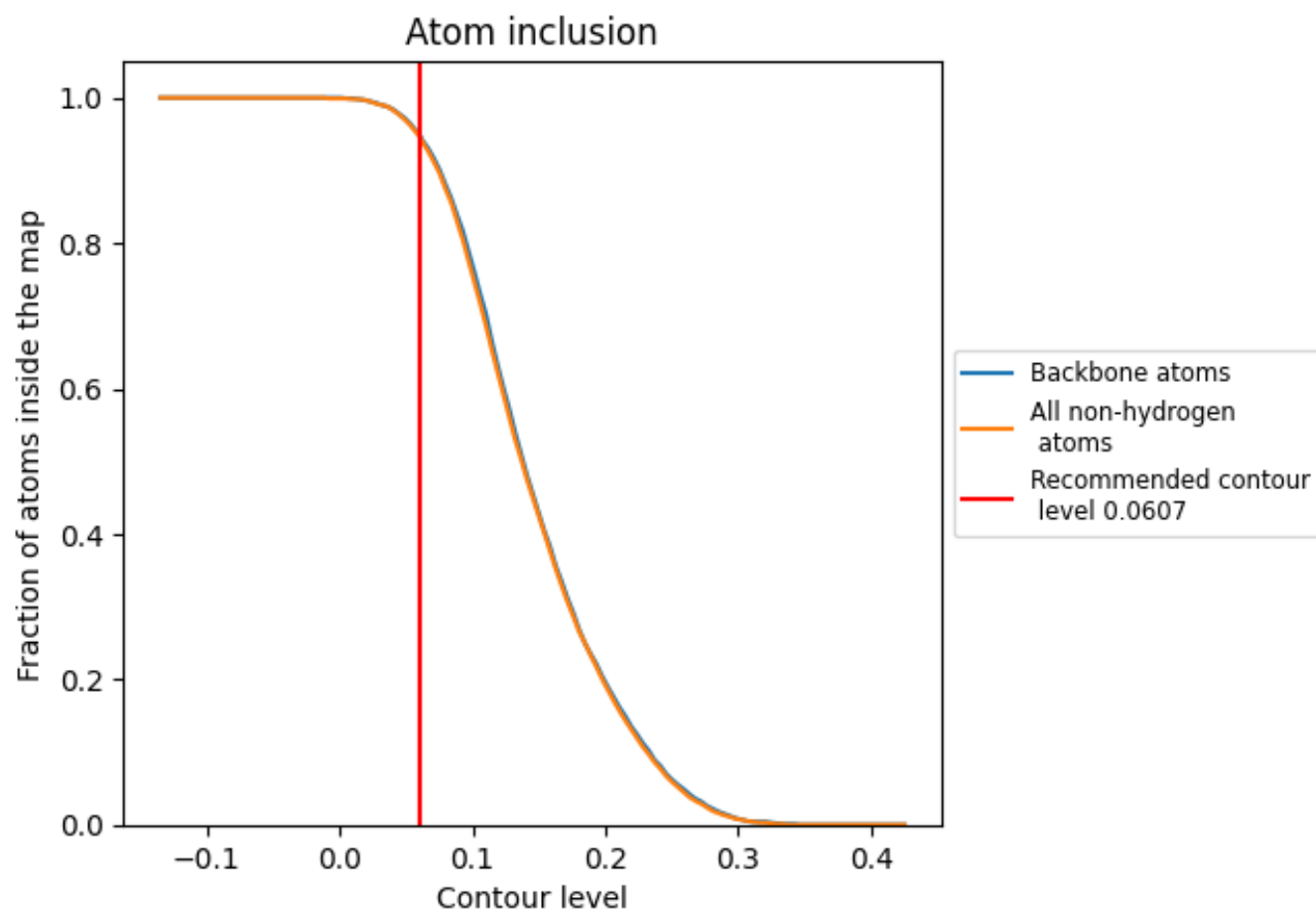
## 9.3 Atom inclusion mapped to coordinate model [i](#)



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 (0.0607).





































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 94% 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 (0.0607) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9450	 0.2480
A	 0.9190	 0.2350
B	 0.9670	 0.2570
C	 0.9520	 0.2460
D	 0.9490	 0.2530
E	 1.0000	 0.3320
F	 0.9400	 0.1930
G	 0.8570	 0.2910
H	 0.9390	 0.1380
I	 0.8930	 0.2210
J	 0.9600	 0.2600
K	 1.0000	 0.3920
L	 1.0000	 0.3700
M	 1.0000	 0.3540
N	 0.9740	 0.2790
O	 0.9640	 0.2820
P	 0.9490	 0.3520
Q	 0.9640	 0.3040
R	 0.8720	 0.2020
S	 0.9740	 0.3720
T	 1.0000	 0.2710
U	 1.0000	 0.3120
V	 1.0000	 0.2850
W	 0.9740	 0.2400
X	 0.9740	 0.2470
Y	 0.9740	 0.2210
Z	 0.9400	 0.3100
a	 0.9840	 0.2990
b	 0.9740	 0.2760
c	 0.9800	 0.3100
d	 1.0000	 0.2700
e	 1.0000	 0.3960
f	 0.7860	 0.1960

