



wwPDB EM Validation Summary Report ⓘ

Mar 9, 2026 – 04:10 PM UTC

PDB ID : 9BD9 / pdb_00009bd9
EMDB ID : EMD-44447
Title : SARS CoV-2 full-length WT spike protein, 1RBD-up conformation (SPIKE-WT)
Authors : Singh, S.; Hasan, S.S.
Deposited on : 2024-04-11
Resolution : 2.95 Å (reported)
Based on initial model : 7KRR

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

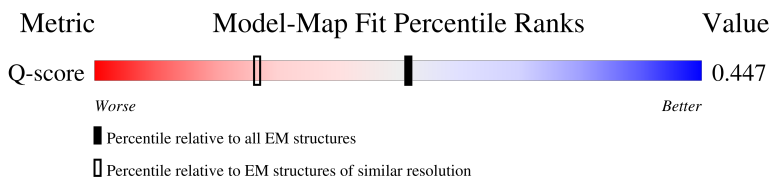
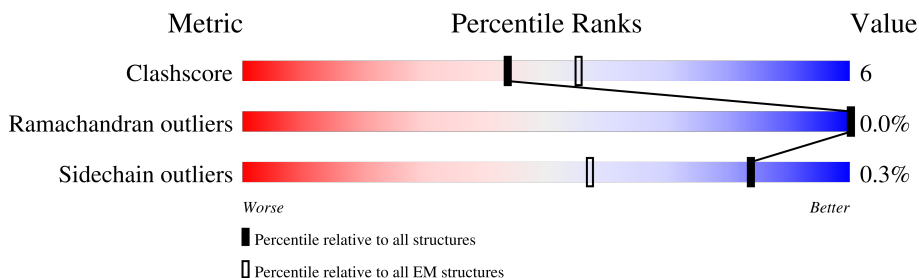
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13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

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.95 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13114 (2.45 - 3.45)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1312	71% 8% 21%
1	B	1312	72% 10% 18%
1	C	1312	74% 9% 17%
2	D	4	25% 75%

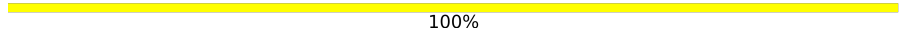



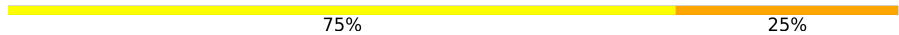




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Mol	Chain	Length	Quality of chain
2	K	4	25% 75%
2	e	4	25% 75%
3	E	2	50% 50%
3	F	2	100%
3	I	2	100%
3	O	2	100%
3	Q	2	100%
3	R	2	100%
3	S	2	50% 50%
3	a	2	100%
3	b	2	100%
3	c	2	100%
3	d	2	50% 50%
3	f	2	100%
4	G	3	33% 67%
4	H	3	33% 67%
4	J	3	33% 67%
4	N	3	67% 33%
4	P	3	33% 67%
4	T	3	33% 67%
4	U	3	33% 67%
4	g	3	33% 67%
5	L	3	33% 67%
5	X	3	33% 67%
5	i	3	33% 67%

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Mol	Chain	Length	Quality of chain
6	M	5	 100%
7	V	4	 25% 75%
7	W	4	 50% 50%
7	h	4	 25% 75%
7	j	4	 75% 25%
8	Y	5	 20% 80%
9	Z	6	 100%
10	k	2	 50% 50%
11	l	2	 100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
10	NAG	k	1	X	-	-	-
11	NAG	l	1	-	-	X	-
12	NAG	A	1404	X	-	-	-
3	NAG	F	1	X	-	-	-
7	MAN	W	4	-	-	X	-
7	NAG	h	1	X	-	-	-
8	BMA	Y	3	-	-	X	-
8	MAN	Y	5	-	-	X	-
9	NAG	Z	2	-	-	X	-
9	FUC	Z	6	-	-	X	-

2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 51610 atoms, of which 25066 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
			Total	C	H	N	O	S		
1	A	1032	15871	5148	7822	1337	1528	36	0	0
1	B	1080	16544	5366	8140	1399	1601	38	0	0
1	C	1091	16767	5435	8254	1418	1622	38	0	0

There are 135 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
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
B	-1	GLY	-	insertion	UNP P0DTC2
B	0	SER	-	insertion	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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
C	-6	GLY	-	insertion	UNP P0DTC2
C	-5	GLY	-	insertion	UNP P0DTC2
C	-4	GLY	-	insertion	UNP P0DTC2
C	-3	SER	-	insertion	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
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

- Molecule 2 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	
2	D	4	Total	C	H	N	O	0	0
			93	28	43	2	20		
2	K	4	Total	C	H	N	O	0	0
			93	28	43	2	20		

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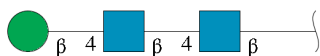
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
2	e	4	93	28	43	2	20	0	0

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



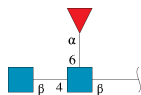
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
3	E	2	40	16	12	2	10	0	0
3	F	2	39	16	11	2	10	0	0
3	I	2	53	16	25	2	10	0	0
3	O	2	40	16	12	2	10	0	0
3	Q	2	40	16	12	2	10	0	0
3	R	2	53	16	25	2	10	0	0
3	S	2	28	16	2	2	10	0	0
3	a	2	53	16	25	2	10	0	0
3	b	2	53	16	25	2	10	0	0
3	c	2	40	16	12	2	10	0	0
3	d	2	53	16	25	2	10	0	0
3	f	2	53	16	25	2	10	0	0

- Molecule 4 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.



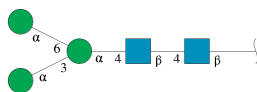
Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	3	Total	C	H	N	O	0	0
			73	22	34	2	15		
4	H	3	Total	C	H	N	O	0	0
			63	22	24	2	15		
4	J	3	Total	C	H	N	O	0	0
			73	22	34	2	15		
4	N	3	Total	C	N	O		0	0
			39	22	2	15			
4	P	3	Total	C	H	N	O	0	0
			73	22	34	2	15		
4	T	3	Total	C	N	O		0	0
			39	22	2	15			
4	U	3	Total	C	N	O		0	0
			39	22	2	15			
4	g	3	Total	C	H	N	O	0	0
			73	22	34	2	15		

- Molecule 5 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
5	L	3	Total	C	H	N	O	0	0
			59	22	21	2	14		
5	X	3	Total	C	N	O		0	0
			38	22	2	14			
5	i	3	Total	C	H	N	O	0	0
			72	22	34	2	14		

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



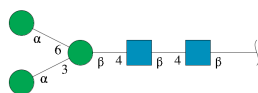
Mol	Chain	Residues	Atoms					AltConf	Trace
6	M	5	Total	C	H	N	O	0	0
			113	34	52	2	25		

- Molecule 7 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.



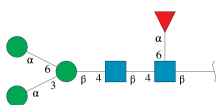
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
7	V	4	50	28	2	20	0	0
7	W	4	50	28	2	20	0	0
7	h	4	93	28	43	20	0	0
7	j	4	74	28	24	20	0	0

- 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
			Total	C	N	O		
8	Y	5	61	34	2	25	0	0

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



Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
9	Z	6	71	40	2	29	0	0

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



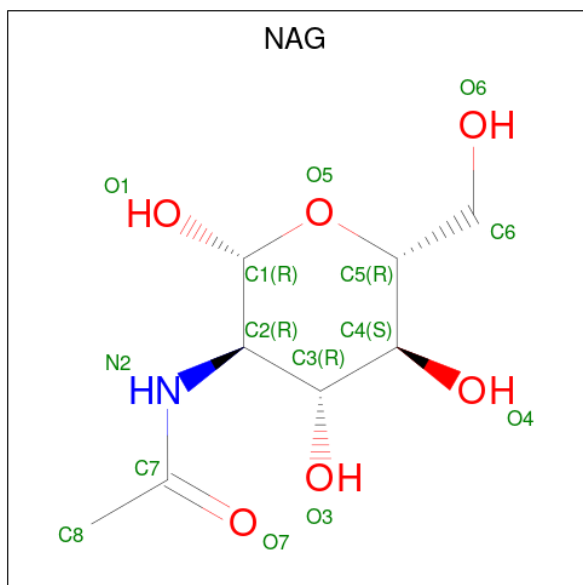
Mol	Chain	Residues	Atoms				AltConf	Trace	
			Total	C	H	N			O
10	k	2	34	14	10	1	9	0	0

- Molecule 11 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
11	l	2	25	14	1	10	0	0

- Molecule 12 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

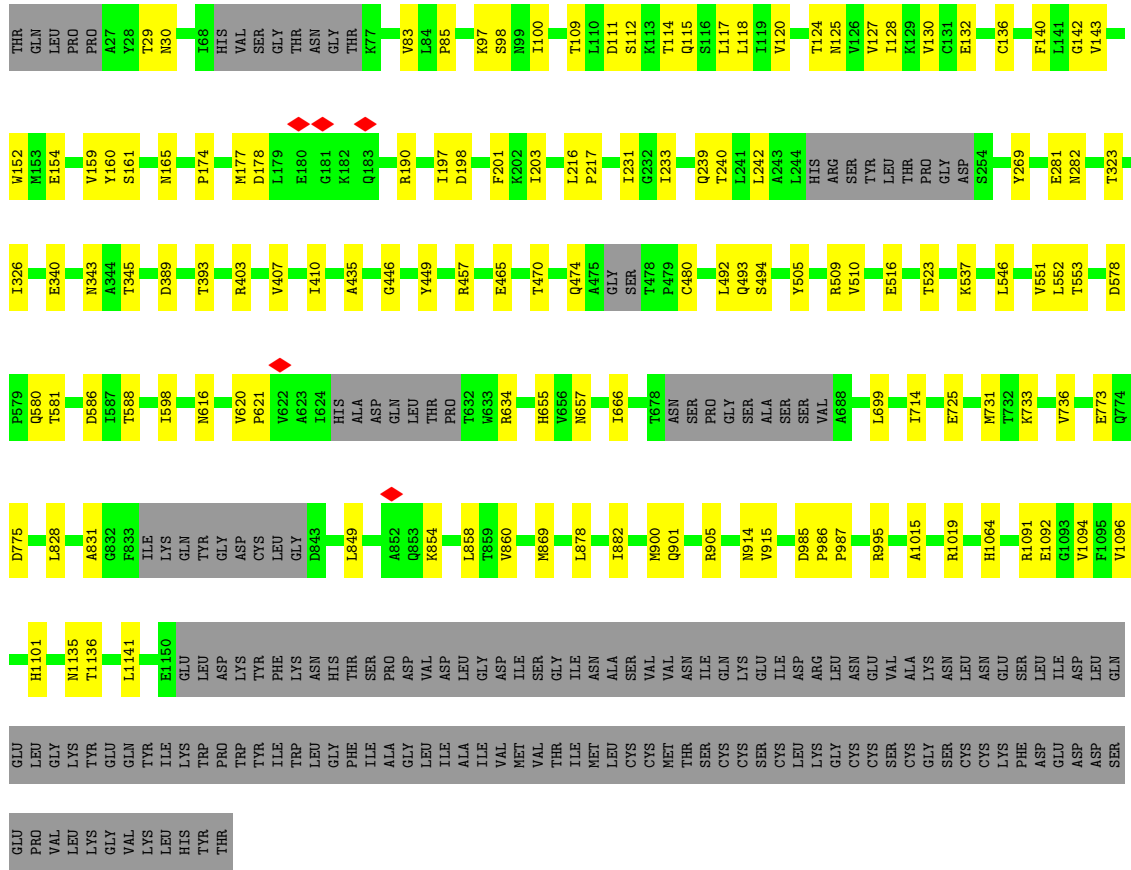


Mol	Chain	Residues	Atoms					AltConf
			Total	C	H	N	O	
12	A	1	27	8	13	1	5	0
12	A	1	27	8	13	1	5	0

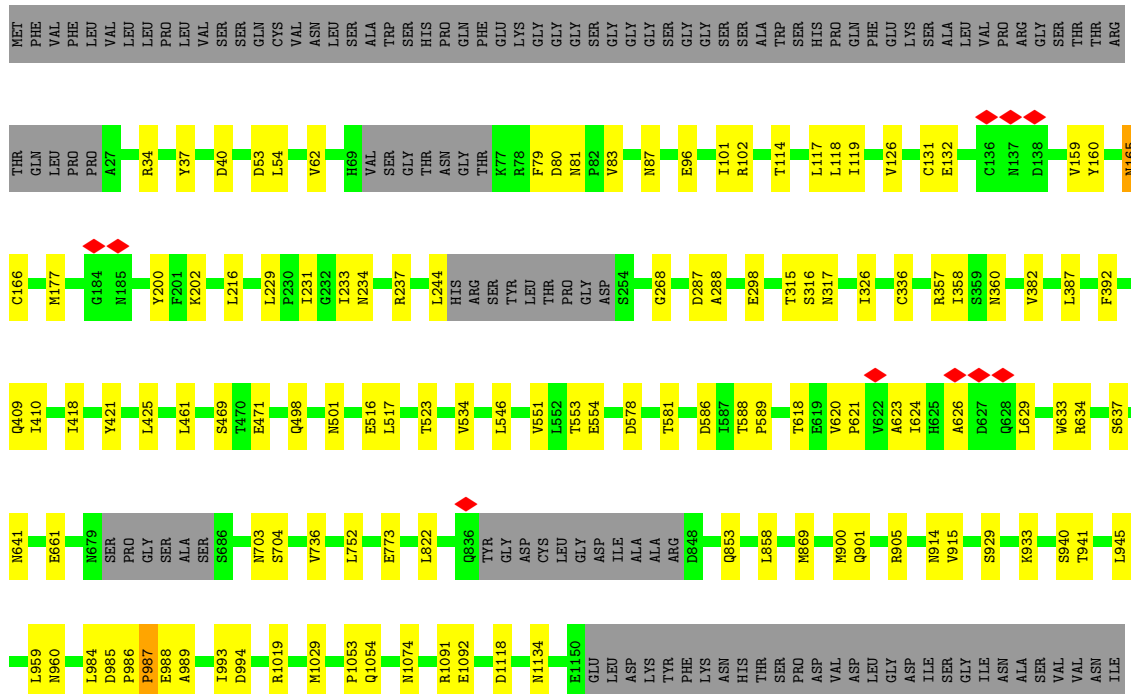
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Mol	Chain	Residues	Atoms					AltConf
12	A	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	A	1	Total	C	H	N	O	0
			26	8	12	1	5	
12	A	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	A	1	Total	C	N	O		0
			14	8	1	5		
12	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	B	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	B	1	Total	C	N	O		0
			14	8	1	5		
12	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	C	1	Total	C	H	N	O	0
			27	8	13	1	5	
12	C	1	Total	C	N	O		0
			14	8	1	5		



● Molecule 1: Spike glycoprotein



GLN
LYS
GLU
ILE
ASP
ARG
LEU
ASN
GLU
VAL
SER
ALA
LYS
ASN
LEU
ASN
GLU
SER
PHE
SER
LEU
ILE
ASP
LEU
GLN
SER
GLU
LEU
PRO
VAL
GLY
LYS
TYR
GLU
GLN
TYR
ILE
LYS
TRP
TRP
PRO
TYR
TRP
ILE
TRP
LEU
GLY
PHE
ILE
ALA
GLY
LEU
ILE
ALA
ILE
VAL
MET
VAL
THR
THR
MET
LEU
CYS
CYS
MET
THR
SER

CYS
CYS
SER
CYS
LEU
LYS
GLY
CYS
CYS
SER
CYS
ALA
GLY
SER
CYS
CYS
LYS
PHE
ASP
GLU
ASP
SER
GLU
PRO
VAL
LEU
LYS
GLY
VAL
LYS
HIS
TYR
THR

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

Chain D:  25% 75%

MAG1
MAG2
BMA3
MAM4

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

Chain K:  25% 75%

MAG1
MAG2
BMA3
MAM4

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

Chain e:  25% 75%

MAG1
MAG2
BMA3
MAM4

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

Chain E:  50% 50%

MAG1
MAG2

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

Chain F:  100%

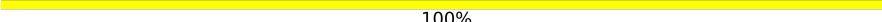
MAG1
MAG2

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

Chain I:  100%

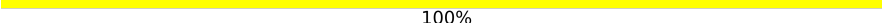
MAG1
MAG2

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

Chain O:  100%

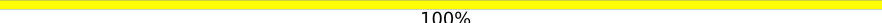
MAG1
MAG2

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

Chain Q:  100%

MAG1
MAG2

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

Chain R:  100%

MAG1
MAG2

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

Chain S:  50%

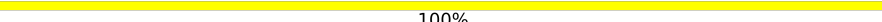
MAG1
MAG2

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

Chain a:  100%

MAG1
MAG2

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

Chain b:  100%


MAG1
MAG2

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

Chain c:  100%

MAG1
MAG2

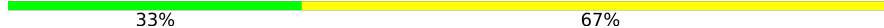
- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain d:  50% 50%MAG1
MAG2

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

Chain f:  100%MAG1
MAG2


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

Chain G:  33% 67%MAG1
MAG2
BMA3

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

Chain H:  33% 67%MAG1
MAG2
BMA3

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

Chain J:  33% 67%MAG1
MAG2
BMA3

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

Chain N:  67% 33%MAG1
MAG2
BMA3

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

Chain P:  33% 67%



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

Chain T:  33% 67%

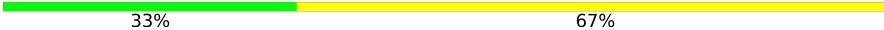


- Molecule 4: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain U:  33% 67%



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

Chain g:  33% 67%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  33% 67%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain X:  33% 67%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain i:  33% 67%



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

Chain M: 100%



- Molecule 7: 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: 25% 75%



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

Chain W: 50% 50%



- Molecule 7: 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 h: 25% 75%



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

Chain j: 75% 25%



- 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 Y: 20% 80%



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

Chain Z:  100%

MAG1
MAG2
BMA3
MAN4
MAN5
FUC6

- Molecule 10: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain k:  50%  50%

MAG1
FUC2

- Molecule 11: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain l:  100%

MAG1
BMA2

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	117450	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOCONTINUUM (6k x 4k)	Depositor
Maximum map value	0.924	Depositor
Minimum map value	-0.303	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.024	Depositor
Recommended contour level	0.08	Depositor
Map size (\AA)	350.01, 350.01, 350.01	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.1667, 1.1667, 1.1667	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.17	0/8231	0.26	0/11196
1	B	0.17	0/8596	0.25	0/11696
1	C	0.17	0/8711	0.25	0/11854
All	All	0.17	0/25538	0.25	0/34746

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	8049	7822	7848	76	0
1	B	8404	8140	8160	92	0
1	C	8513	8254	8280	91	0
2	D	50	43	43	0	0
2	K	50	43	43	0	0
2	e	50	43	43	0	0
3	E	28	12	25	1	0
3	F	28	11	24	3	0
3	I	28	25	25	1	0
3	O	28	12	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	Q	28	12	24	0	0
3	R	28	25	25	0	0
3	S	28	0	25	5	0
3	a	28	25	25	1	0
3	b	28	25	25	0	0
3	c	28	12	24	3	0
3	d	28	25	25	2	0
3	f	28	25	25	2	0
4	G	39	34	34	0	0
4	H	39	24	34	2	0
4	J	39	34	34	0	0
4	N	39	0	34	2	0
4	P	39	34	34	2	0
4	T	39	0	34	3	0
4	U	39	0	34	6	0
4	g	39	34	34	0	0
5	L	38	21	34	3	0
5	X	38	0	34	4	0
5	i	38	34	34	1	0
6	M	61	52	52	0	0
7	V	50	0	43	4	0
7	W	50	0	43	7	0
7	h	50	43	43	2	0
7	j	50	24	43	3	0
8	Y	61	0	52	20	0
9	Z	71	0	61	16	0
10	k	24	10	21	6	0
11	l	25	0	22	7	0
12	A	84	64	77	6	0
12	B	70	52	65	2	0
12	C	70	52	65	0	0
All	All	26544	25066	25675	337	0

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

The worst 5 of 337 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)
11:l:1:NAG:H5	11:l:2:BMA:H2	1.18	1.11
7:W:3:BMA:H5	7:W:4:MAN:H3	1.35	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:Y:3:BMA:H5	8:Y:5:MAN:H5	1.42	1.02
4:U:1:NAG:HN2	4:U:2:NAG:H2	1.26	0.96
7:j:3:BMA:H62	7:j:4:MAN:H3	1.49	0.94

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	A	1014/1312 (77%)	964 (95%)	50 (5%)	0	100	100
1	B	1066/1312 (81%)	1007 (94%)	59 (6%)	0	100	100
1	C	1081/1312 (82%)	1017 (94%)	63 (6%)	1 (0%)	48	72
All	All	3161/3936 (80%)	2988 (94%)	172 (5%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	987	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	895/1137 (79%)	889 (99%)	6 (1%)	76	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	930/1137 (82%)	930 (100%)	0	100	100
1	C	946/1137 (83%)	945 (100%)	1 (0%)	88	95
All	All	2771/3411 (81%)	2764 (100%)	7 (0%)	84	92

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	855	PHE
1	A	923	ILE
1	C	165	ASN
1	A	1002	GLN
1	A	657	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	762	GLN
1	C	1002	GLN
1	A	658	ASN
1	A	762	GLN
1	A	919	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

105 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	2,1	14,14,15	2.02	5 (35%)	17,19,21	1.18	1 (5%)
2	NAG	D	2	2	14,14,15	1.99	4 (28%)	17,19,21	1.01	2 (11%)
2	BMA	D	3	2	11,11,12	0.52	0	15,15,17	0.61	0
2	MAN	D	4	2	11,11,12	0.55	0	15,15,17	1.03	2 (13%)
3	NAG	E	1	3,1	14,14,15	1.95	4 (28%)	17,19,21	2.61	7 (41%)
3	NAG	E	2	3	14,14,15	2.05	4 (28%)	17,19,21	1.09	1 (5%)
3	NAG	F	1	3,1	14,14,15	1.94	4 (28%)	17,19,21	2.74	8 (47%)
3	NAG	F	2	3	14,14,15	2.01	4 (28%)	17,19,21	1.39	2 (11%)
4	NAG	G	1	4,1	14,14,15	2.06	5 (35%)	17,19,21	1.14	2 (11%)
4	NAG	G	2	4	14,14,15	2.01	4 (28%)	17,19,21	1.18	2 (11%)
4	BMA	G	3	4	11,11,12	0.74	0	15,15,17	0.73	0
4	NAG	H	1	4,1	14,14,15	1.98	5 (35%)	17,19,21	1.66	4 (23%)
4	NAG	H	2	4	14,14,15	1.98	4 (28%)	17,19,21	1.32	2 (11%)
4	BMA	H	3	4	11,11,12	0.54	0	15,15,17	0.65	0
3	NAG	I	1	3,1	14,14,15	2.06	5 (35%)	17,19,21	1.16	1 (5%)
3	NAG	I	2	3	14,14,15	2.02	4 (28%)	17,19,21	0.87	1 (5%)
4	NAG	J	1	4,1	14,14,15	1.90	3 (21%)	17,19,21	1.21	2 (11%)
4	NAG	J	2	4	14,14,15	1.92	3 (21%)	17,19,21	1.07	2 (11%)
4	BMA	J	3	4	11,11,12	0.50	0	15,15,17	0.64	0
2	NAG	K	1	2,1	14,14,15	1.89	3 (21%)	17,19,21	1.11	2 (11%)
2	NAG	K	2	2	14,14,15	2.04	4 (28%)	17,19,21	0.84	0
2	BMA	K	3	2	11,11,12	0.56	0	15,15,17	0.67	0
2	MAN	K	4	2	11,11,12	1.19	1 (9%)	15,15,17	1.87	2 (13%)
5	NAG	L	1	5,1	14,14,15	1.89	3 (21%)	17,19,21	1.55	5 (29%)
5	NAG	L	2	5	14,14,15	1.99	4 (28%)	17,19,21	1.32	2 (11%)
5	FUC	L	3	5	10,10,11	0.56	0	14,14,16	0.69	0
6	NAG	M	1	6,1	14,14,15	1.93	4 (28%)	17,19,21	1.11	2 (11%)
6	NAG	M	2	6	14,14,15	1.86	3 (21%)	17,19,21	1.51	4 (23%)
6	MAN	M	3	6	11,11,12	0.99	1 (9%)	15,15,17	1.01	1 (6%)
6	MAN	M	4	6	11,11,12	0.58	0	15,15,17	1.02	2 (13%)
6	MAN	M	5	6	11,11,12	0.65	0	15,15,17	0.89	1 (6%)
4	NAG	N	1	4,1	14,14,15	2.12	4 (28%)	17,19,21	2.03	5 (29%)
4	NAG	N	2	4	14,14,15	2.09	4 (28%)	17,19,21	1.14	1 (5%)
4	BMA	N	3	4	11,11,12	0.76	0	15,15,17	0.75	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	O	1	3,1	14,14,15	1.96	4 (28%)	17,19,21	1.02	1 (5%)
3	NAG	O	2	3	14,14,15	2.04	4 (28%)	17,19,21	1.14	2 (11%)
4	NAG	P	1	4,1	14,14,15	1.97	4 (28%)	17,19,21	1.23	2 (11%)
4	NAG	P	2	4	14,14,15	2.02	4 (28%)	17,19,21	0.96	1 (5%)
4	BMA	P	3	4	11,11,12	0.52	0	15,15,17	0.67	0
3	NAG	Q	1	3,1	14,14,15	1.92	3 (21%)	17,19,21	1.62	3 (17%)
3	NAG	Q	2	3	14,14,15	1.98	4 (28%)	17,19,21	1.11	2 (11%)
3	NAG	R	1	3,1	14,14,15	1.96	4 (28%)	17,19,21	1.18	2 (11%)
3	NAG	R	2	3	14,14,15	2.01	4 (28%)	17,19,21	1.04	2 (11%)
3	NAG	S	1	3,1	14,14,15	2.42	4 (28%)	17,19,21	1.99	5 (29%)
3	NAG	S	2	3	14,14,15	2.13	4 (28%)	17,19,21	1.26	2 (11%)
4	NAG	T	1	4,1	14,14,15	1.94	3 (21%)	17,19,21	1.20	2 (11%)
4	NAG	T	2	4	14,14,15	2.01	4 (28%)	17,19,21	1.00	1 (5%)
4	BMA	T	3	4	11,11,12	0.54	0	15,15,17	0.71	0
4	NAG	U	1	4,1	14,14,15	2.04	4 (28%)	17,19,21	1.63	3 (17%)
4	NAG	U	2	4	14,14,15	1.99	4 (28%)	17,19,21	1.19	2 (11%)
4	BMA	U	3	4	11,11,12	0.51	0	15,15,17	0.66	0
7	NAG	V	1	7,1	14,14,15	1.94	5 (35%)	17,19,21	1.14	2 (11%)
7	NAG	V	2	7	14,14,15	1.98	4 (28%)	17,19,21	1.22	2 (11%)
7	BMA	V	3	7	11,11,12	0.87	0	15,15,17	1.31	2 (13%)
7	MAN	V	4	7	11,11,12	0.56	0	15,15,17	0.95	2 (13%)
7	NAG	W	1	7,1	14,14,15	1.90	3 (21%)	17,19,21	1.17	2 (11%)
7	NAG	W	2	7	14,14,15	1.99	4 (28%)	17,19,21	0.96	1 (5%)
7	BMA	W	3	7	11,11,12	0.59	0	15,15,17	0.67	0
7	MAN	W	4	7	11,11,12	1.08	1 (9%)	15,15,17	1.65	2 (13%)
5	NAG	X	1	5,1	14,14,15	2.05	4 (28%)	17,19,21	1.09	1 (5%)
5	NAG	X	2	5	14,14,15	1.99	4 (28%)	17,19,21	1.12	2 (11%)
5	FUC	X	3	5	10,10,11	0.80	0	14,14,16	0.91	1 (7%)
8	NAG	Y	1	8,1	14,14,15	2.00	5 (35%)	17,19,21	1.25	2 (11%)
8	NAG	Y	2	8	14,14,15	2.04	4 (28%)	17,19,21	1.24	3 (17%)
8	BMA	Y	3	8	11,11,12	0.72	0	15,15,17	0.92	0
8	MAN	Y	4	8	11,11,12	0.66	0	15,15,17	0.91	2 (13%)
8	MAN	Y	5	8	11,11,12	1.03	1 (9%)	15,15,17	1.31	3 (20%)
9	NAG	Z	1	9,1	14,14,15	1.93	4 (28%)	17,19,21	1.31	4 (23%)
9	NAG	Z	2	9	14,14,15	1.94	4 (28%)	17,19,21	1.27	2 (11%)
9	BMA	Z	3	9	11,11,12	0.64	0	15,15,17	0.95	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	MAN	Z	4	9	11,11,12	0.62	0	15,15,17	1.03	2 (13%)
9	MAN	Z	5	9	11,11,12	0.57	0	15,15,17	0.99	2 (13%)
9	FUC	Z	6	9	10,10,11	1.17	1 (10%)	14,14,16	1.63	3 (21%)
3	NAG	a	1	3,1	14,14,15	2.02	4 (28%)	17,19,21	1.10	2 (11%)
3	NAG	a	2	3	14,14,15	1.99	4 (28%)	17,19,21	1.03	1 (5%)
3	NAG	b	1	3,1	14,14,15	1.93	4 (28%)	17,19,21	1.21	2 (11%)
3	NAG	b	2	3	14,14,15	2.03	4 (28%)	17,19,21	1.07	1 (5%)
3	NAG	c	1	3,1	14,14,15	2.07	4 (28%)	17,19,21	2.32	5 (29%)
3	NAG	c	2	3	14,14,15	1.99	4 (28%)	17,19,21	1.16	2 (11%)
3	NAG	d	1	3,1	14,14,15	1.97	3 (21%)	17,19,21	1.40	1 (5%)
3	NAG	d	2	3	14,14,15	1.97	3 (21%)	17,19,21	1.52	2 (11%)
2	NAG	e	1	2,1	14,14,15	1.95	4 (28%)	17,19,21	1.20	2 (11%)
2	NAG	e	2	2	14,14,15	1.93	4 (28%)	17,19,21	1.94	5 (29%)
2	BMA	e	3	2	11,11,12	0.61	0	15,15,17	0.77	0
2	MAN	e	4	2	11,11,12	0.71	0	15,15,17	1.53	2 (13%)
3	NAG	f	1	3,1	14,14,15	1.96	4 (28%)	17,19,21	1.50	2 (11%)
3	NAG	f	2	3	14,14,15	1.97	4 (28%)	17,19,21	1.26	1 (5%)
4	NAG	g	1	4,1	14,14,15	1.93	4 (28%)	17,19,21	1.15	2 (11%)
4	NAG	g	2	4	14,14,15	1.93	4 (28%)	17,19,21	1.15	1 (5%)
4	BMA	g	3	4	11,11,12	0.50	0	15,15,17	0.69	0
7	NAG	h	1	7,1	14,14,15	1.95	3 (21%)	17,19,21	2.50	7 (41%)
7	NAG	h	2	7	14,14,15	2.00	4 (28%)	17,19,21	1.58	2 (11%)
7	BMA	h	3	7	11,11,12	1.02	1 (9%)	15,15,17	0.90	1 (6%)
7	MAN	h	4	7	11,11,12	0.64	0	15,15,17	0.91	1 (6%)
5	NAG	i	1	5,1	14,14,15	1.92	4 (28%)	17,19,21	1.81	4 (23%)
5	NAG	i	2	5	14,14,15	1.99	4 (28%)	17,19,21	1.17	2 (11%)
5	FUC	i	3	5	10,10,11	0.83	0	14,14,16	0.72	0
7	NAG	j	1	7,1	14,14,15	1.91	3 (21%)	17,19,21	1.19	2 (11%)
7	NAG	j	2	7	14,14,15	1.93	4 (28%)	17,19,21	1.20	2 (11%)
7	BMA	j	3	7	11,11,12	0.53	0	15,15,17	0.73	0
7	MAN	j	4	7	11,11,12	1.05	1 (9%)	15,15,17	1.35	3 (20%)
10	NAG	k	1	10,1	14,14,15	2.19	4 (28%)	17,19,21	3.05	11 (64%)
10	FUC	k	2	10	10,10,11	0.62	0	14,14,16	0.81	0
11	NAG	l	1	11	14,14,15	2.11	5 (35%)	17,19,21	1.76	3 (17%)
11	BMA	l	2	11	11,11,12	0.60	0	15,15,17	1.00	1 (6%)

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	2,1	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1
2	BMA	D	3	2	-	0/2/19/22	0/1/1/1
2	MAN	D	4	2	-	2/2/19/22	0/1/1/1
3	NAG	E	1	3,1	-	5/6/23/26	0/1/1/1
3	NAG	E	2	3	-	1/6/23/26	0/1/1/1
3	NAG	F	1	3,1	1/1/6/7	1/6/23/26	0/1/1/1
3	NAG	F	2	3	-	4/6/23/26	0/1/1/1
4	NAG	G	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
4	BMA	G	3	4	-	0/2/19/22	0/1/1/1
4	NAG	H	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	H	2	4	-	1/6/23/26	0/1/1/1
4	BMA	H	3	4	-	1/2/19/22	0/1/1/1
3	NAG	I	1	3,1	-	3/6/23/26	0/1/1/1
3	NAG	I	2	3	-	4/6/23/26	0/1/1/1
4	NAG	J	1	4,1	-	1/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1
4	BMA	J	3	4	-	0/2/19/22	0/1/1/1
2	NAG	K	1	2,1	-	1/6/23/26	0/1/1/1
2	NAG	K	2	2	-	1/6/23/26	0/1/1/1
2	BMA	K	3	2	-	0/2/19/22	0/1/1/1
2	MAN	K	4	2	-	2/2/19/22	0/1/1/1
5	NAG	L	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	L	2	5	-	0/6/23/26	0/1/1/1
5	FUC	L	3	5	-	-	0/1/1/1
6	NAG	M	1	6,1	-	1/6/23/26	0/1/1/1
6	NAG	M	2	6	-	1/6/23/26	0/1/1/1
6	MAN	M	3	6	-	2/2/19/22	1/1/1/1
6	MAN	M	4	6	-	0/2/19/22	0/1/1/1
6	MAN	M	5	6	-	2/2/19/22	0/1/1/1
4	NAG	N	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	N	2	4	-	0/6/23/26	0/1/1/1
4	BMA	N	3	4	-	1/2/19/22	0/1/1/1
3	NAG	O	1	3,1	-	4/6/23/26	0/1/1/1

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	a	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	a	2	3	-	2/6/23/26	0/1/1/1
3	NAG	b	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	b	2	3	-	1/6/23/26	0/1/1/1
3	NAG	c	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	c	2	3	-	2/6/23/26	0/1/1/1
3	NAG	d	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	d	2	3	-	2/6/23/26	0/1/1/1
2	NAG	e	1	2,1	-	1/6/23/26	0/1/1/1
2	NAG	e	2	2	-	2/6/23/26	0/1/1/1
2	BMA	e	3	2	-	1/2/19/22	0/1/1/1
2	MAN	e	4	2	-	0/2/19/22	1/1/1/1
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
4	NAG	g	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	g	2	4	-	2/6/23/26	0/1/1/1
4	BMA	g	3	4	-	1/2/19/22	0/1/1/1
7	NAG	h	1	7,1	1/1/6/7	1/6/23/26	0/1/1/1
7	NAG	h	2	7	-	1/6/23/26	0/1/1/1
7	BMA	h	3	7	-	2/2/19/22	0/1/1/1
7	MAN	h	4	7	-	0/2/19/22	0/1/1/1
5	NAG	i	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	i	2	5	-	2/6/23/26	0/1/1/1
5	FUC	i	3	5	-	-	0/1/1/1
7	NAG	j	1	7,1	-	2/6/23/26	0/1/1/1
7	NAG	j	2	7	-	2/6/23/26	0/1/1/1
7	BMA	j	3	7	-	0/2/19/22	0/1/1/1
7	MAN	j	4	7	-	0/2/19/22	0/1/1/1
10	NAG	k	1	10,1	1/1/6/7	4/6/23/26	0/1/1/1
10	FUC	k	2	10	-	-	0/1/1/1
11	NAG	l	1	11	-	5/6/23/26	0/1/1/1
11	BMA	l	2	11	-	1/2/19/22	0/1/1/1

The worst 5 of 274 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	S	1	NAG	O5-C1	5.38	1.52	1.43
10	k	1	NAG	O5-C1	4.67	1.51	1.43
4	N	2	NAG	O5-C1	4.62	1.51	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	N	1	NAG	O5-C1	4.62	1.51	1.43
5	X	1	NAG	O5-C1	4.57	1.51	1.43

The worst 5 of 207 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	4	MAN	C1-O5-C5	6.30	120.62	112.19
10	k	1	NAG	C3-C4-C5	5.98	121.07	110.23
3	F	1	NAG	C1-O5-C5	-5.95	104.22	112.19
7	W	4	MAN	C1-O5-C5	5.51	119.58	112.19
3	c	1	NAG	O5-C5-C6	5.36	118.10	107.66

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	F	1	NAG	C5
7	h	1	NAG	C3
10	k	1	NAG	C4

5 of 150 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	K	2	NAG	C1-C2-N2-C7
2	e	2	NAG	C1-C2-N2-C7
3	E	1	NAG	C1-C2-N2-C7
3	I	2	NAG	C1-C2-N2-C7
3	O	1	NAG	C1-C2-N2-C7

All (5) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	e	4	MAN	C1-C2-C3-C4-C5-O5
8	Y	4	MAN	C1-C2-C3-C4-C5-O5
7	V	4	MAN	C1-C2-C3-C4-C5-O5
9	Z	4	MAN	C1-C2-C3-C4-C5-O5
6	M	3	MAN	C1-C2-C3-C4-C5-O5

56 monomers are involved in 102 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	N	1	NAG	1	0

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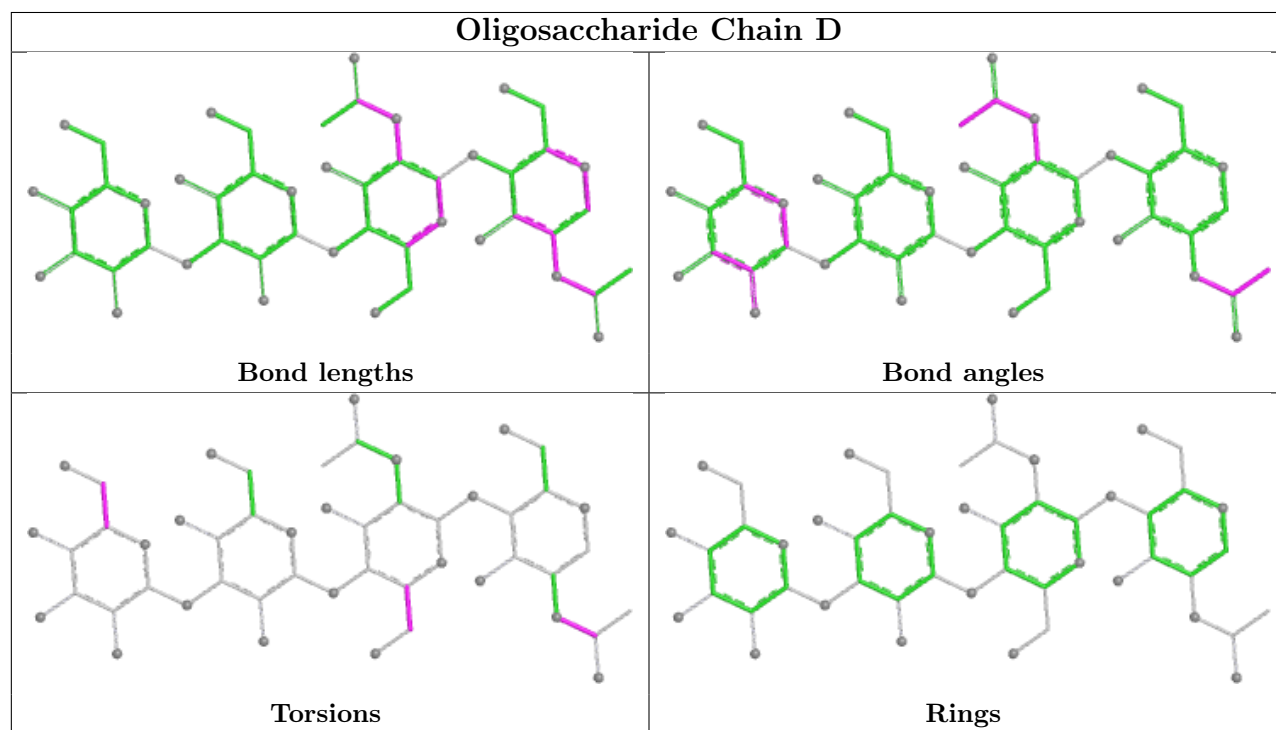
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	1	NAG	1	0
3	F	1	NAG	1	0
10	k	1	NAG	3	0
3	f	1	NAG	1	0
8	Y	4	MAN	4	0
5	X	1	NAG	4	0
10	k	2	FUC	3	0
3	F	2	NAG	3	0
4	T	1	NAG	3	0
5	L	2	NAG	2	0
9	Z	2	NAG	8	0
9	Z	3	BMA	4	0
9	Z	1	NAG	5	0
4	H	2	NAG	2	0
7	h	3	BMA	1	0
11	l	2	BMA	3	0
3	a	2	NAG	1	0
7	W	3	BMA	4	0
4	N	3	BMA	1	0
4	U	2	NAG	3	0
9	Z	5	MAN	5	0
11	l	1	NAG	7	0
7	h	2	NAG	2	0
7	V	3	BMA	3	0
3	d	1	NAG	2	0
4	T	2	NAG	1	0
7	j	3	BMA	3	0
5	L	1	NAG	3	0
4	P	2	NAG	1	0
4	H	3	BMA	1	0
5	i	2	NAG	1	0
3	S	1	NAG	5	0
5	i	1	NAG	1	0
9	Z	6	FUC	7	0
7	j	4	MAN	3	0
7	V	4	MAN	2	0
7	W	2	NAG	3	0
7	W	4	MAN	7	0
5	X	3	FUC	2	0
3	a	1	NAG	1	0
3	E	1	NAG	1	0
4	U	1	NAG	6	0

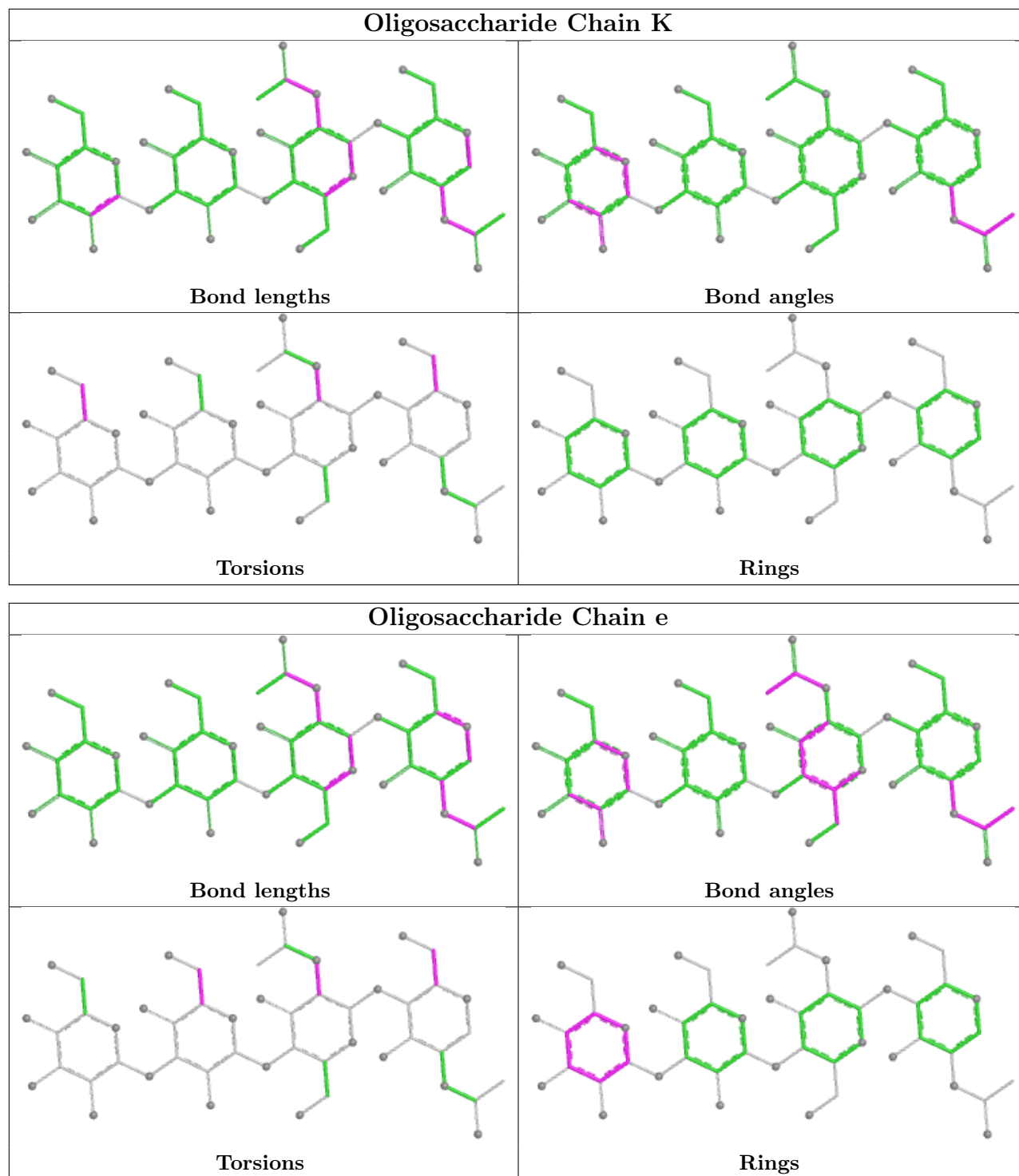
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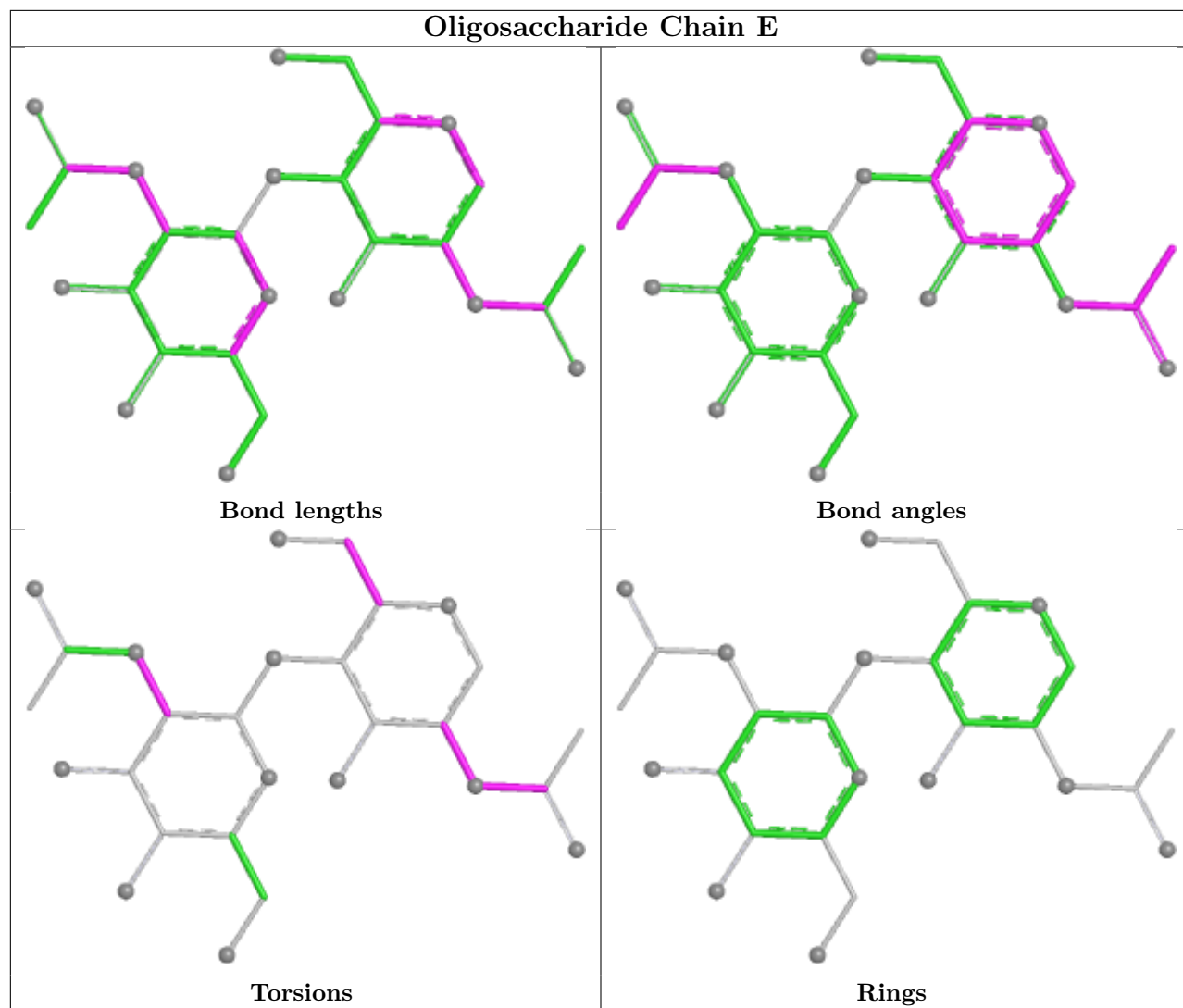
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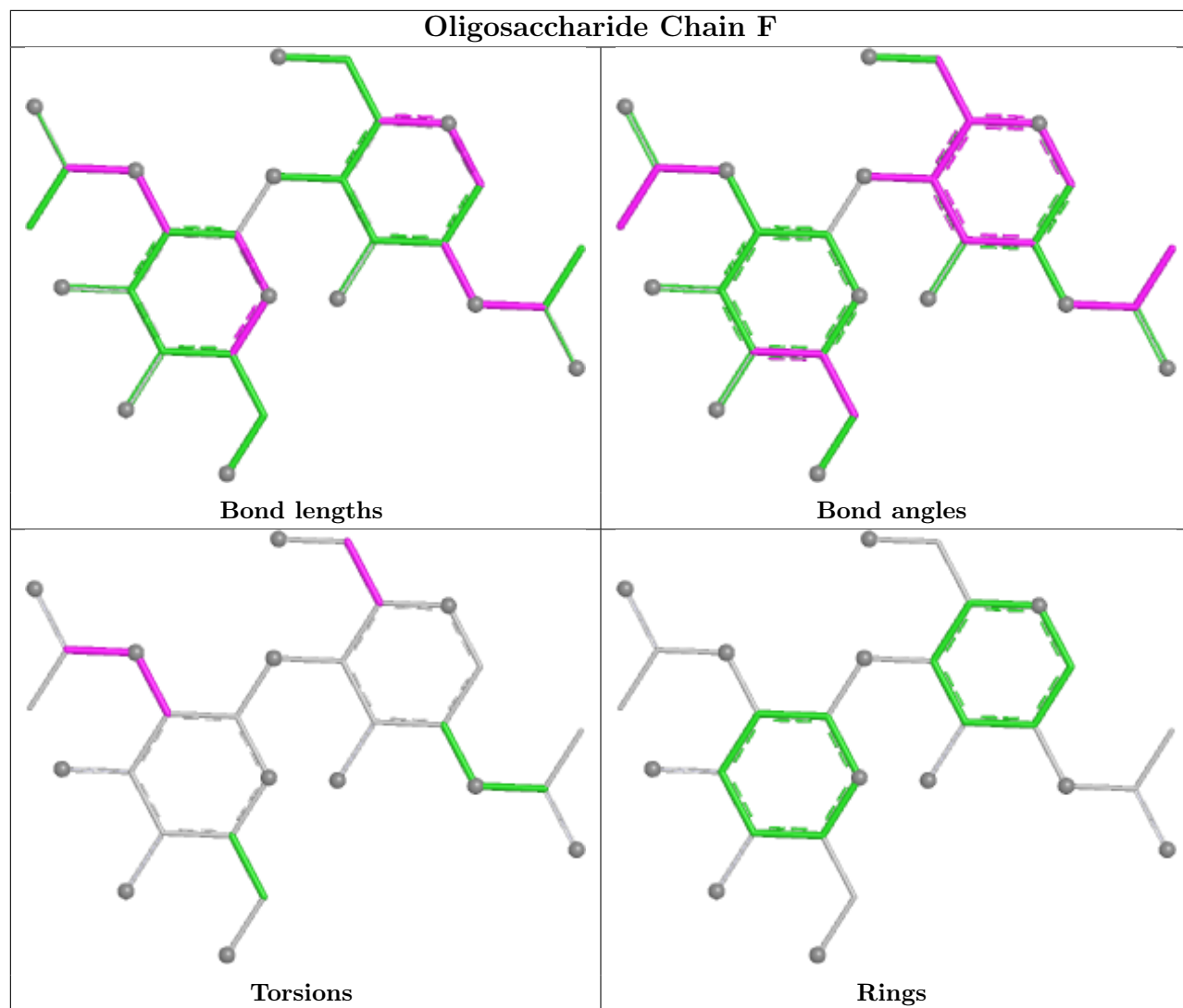
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	Z	4	MAN	3	0
7	h	1	NAG	1	0
3	I	2	NAG	1	0
4	H	1	NAG	1	0
8	Y	5	MAN	9	0
7	V	2	NAG	3	0
4	P	1	NAG	2	0
3	f	2	NAG	1	0
8	Y	3	BMA	14	0
3	c	1	NAG	3	0
8	Y	1	NAG	6	0
8	Y	2	NAG	6	0
3	c	2	NAG	1	0

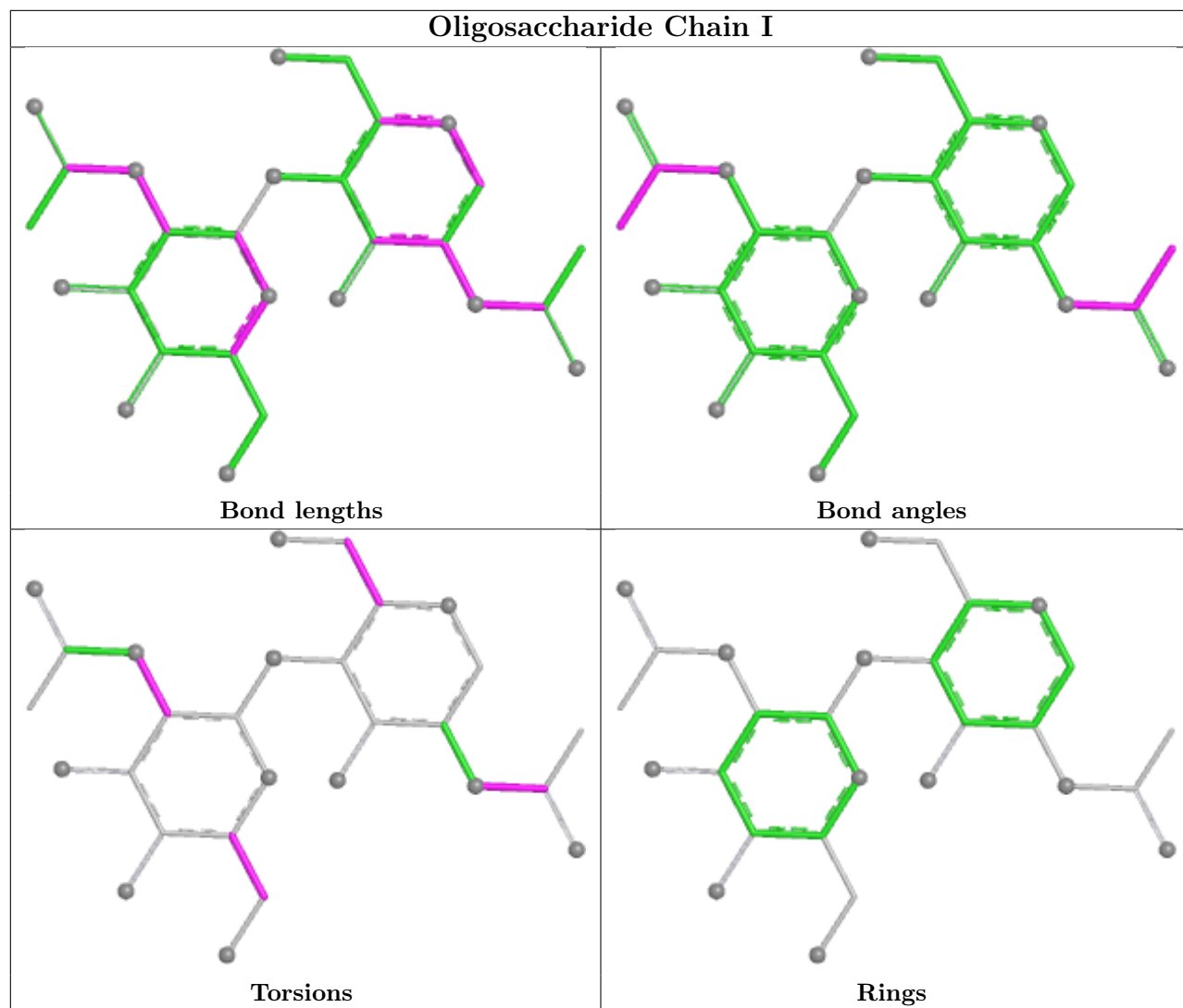
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

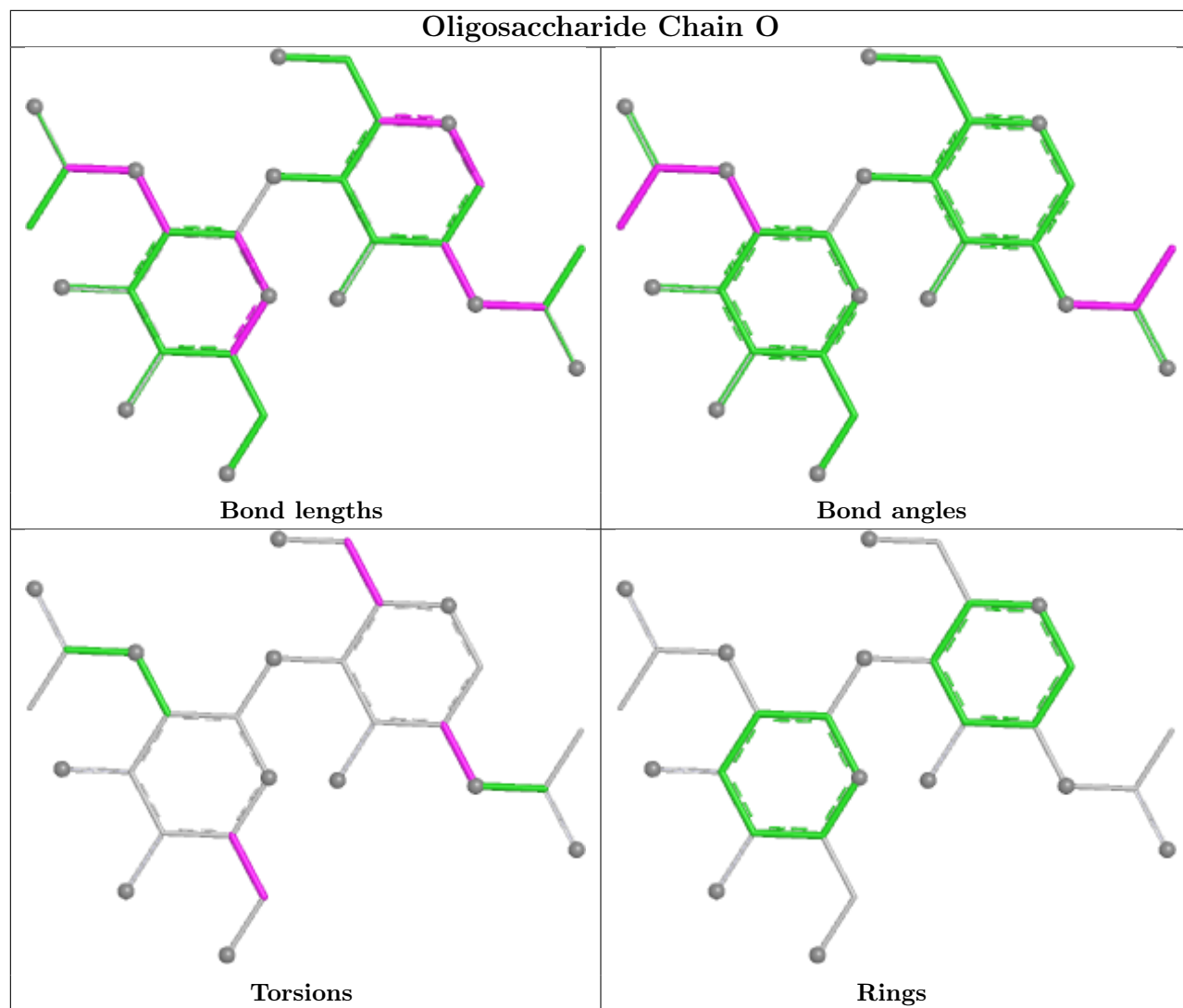


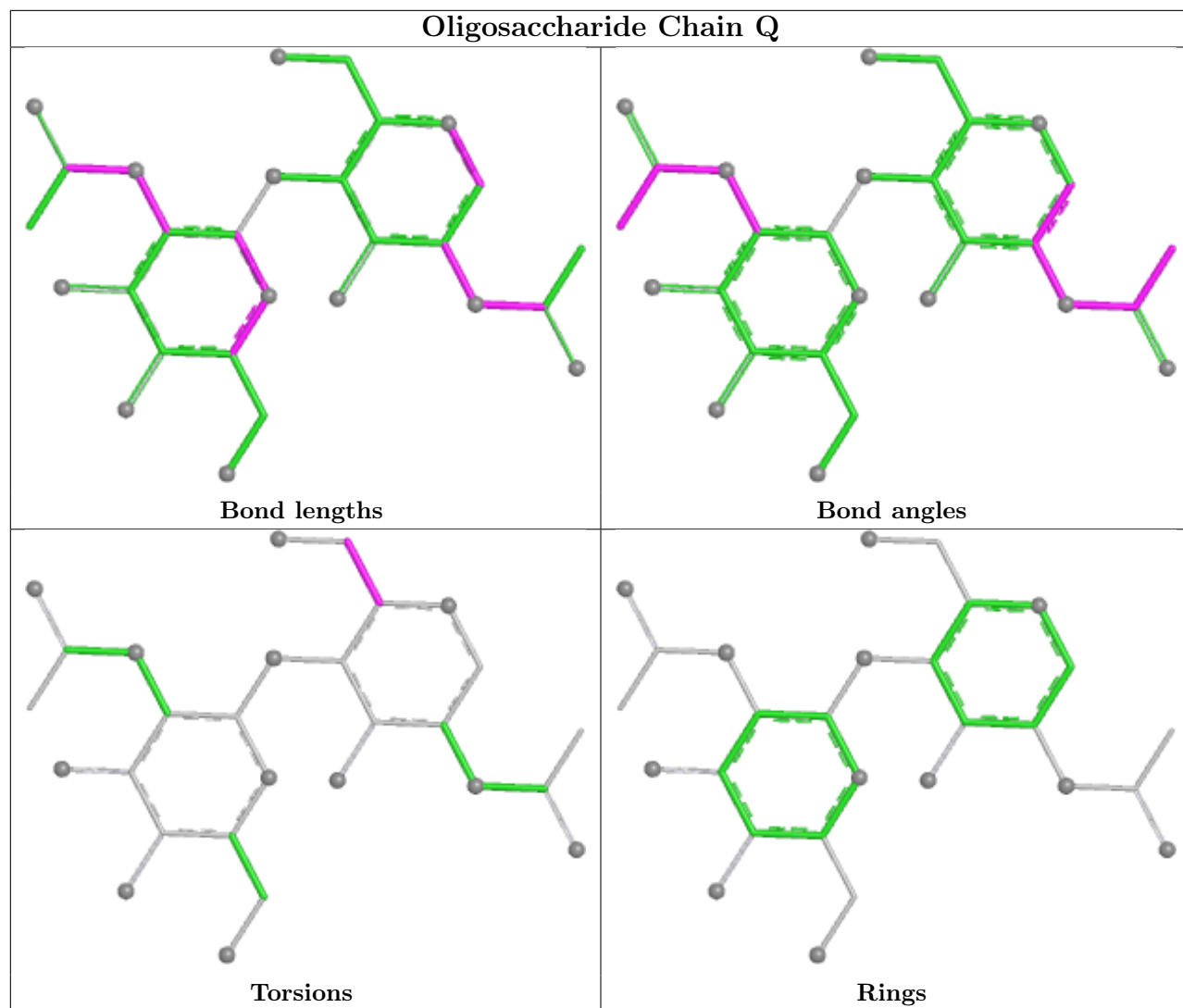


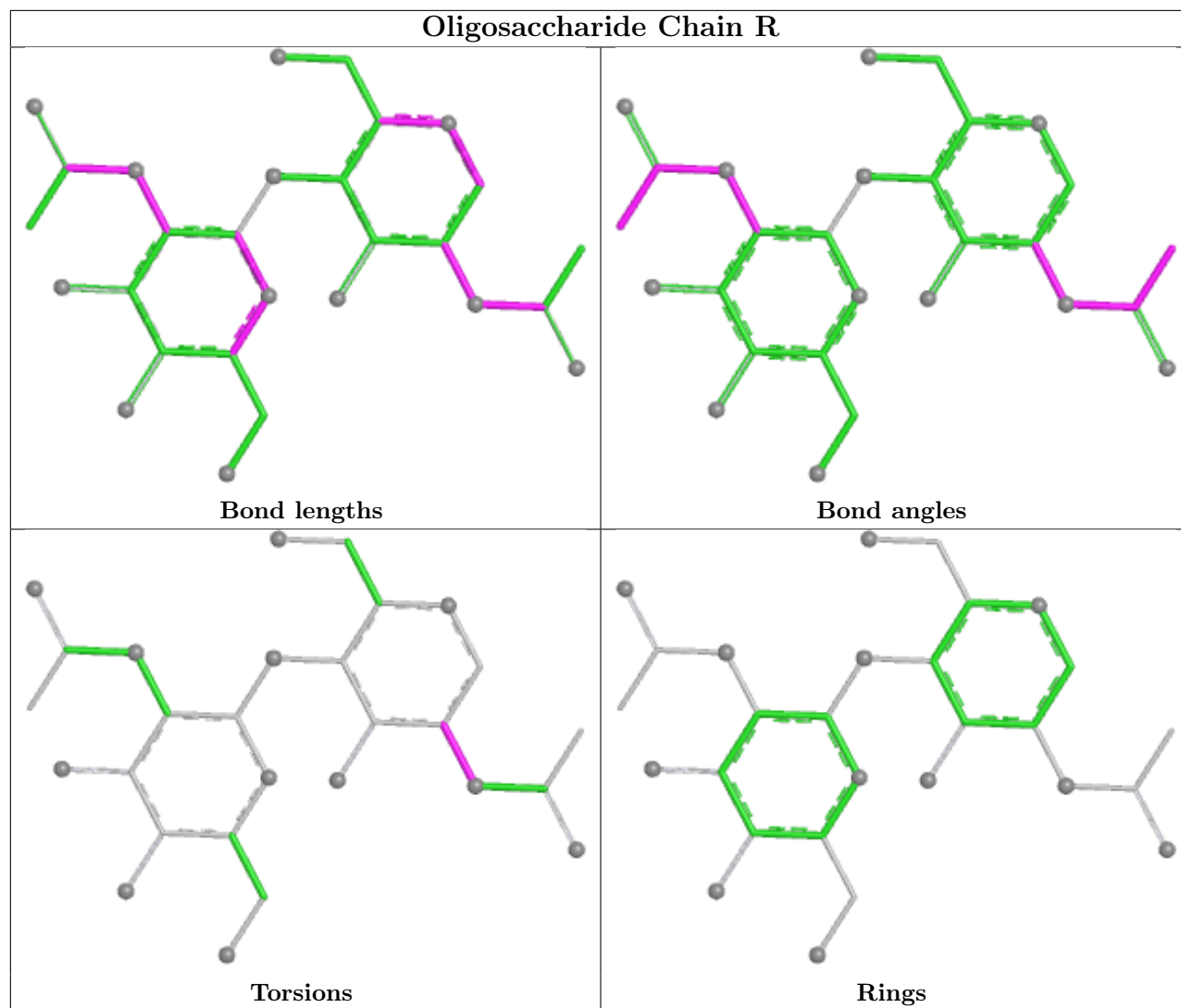


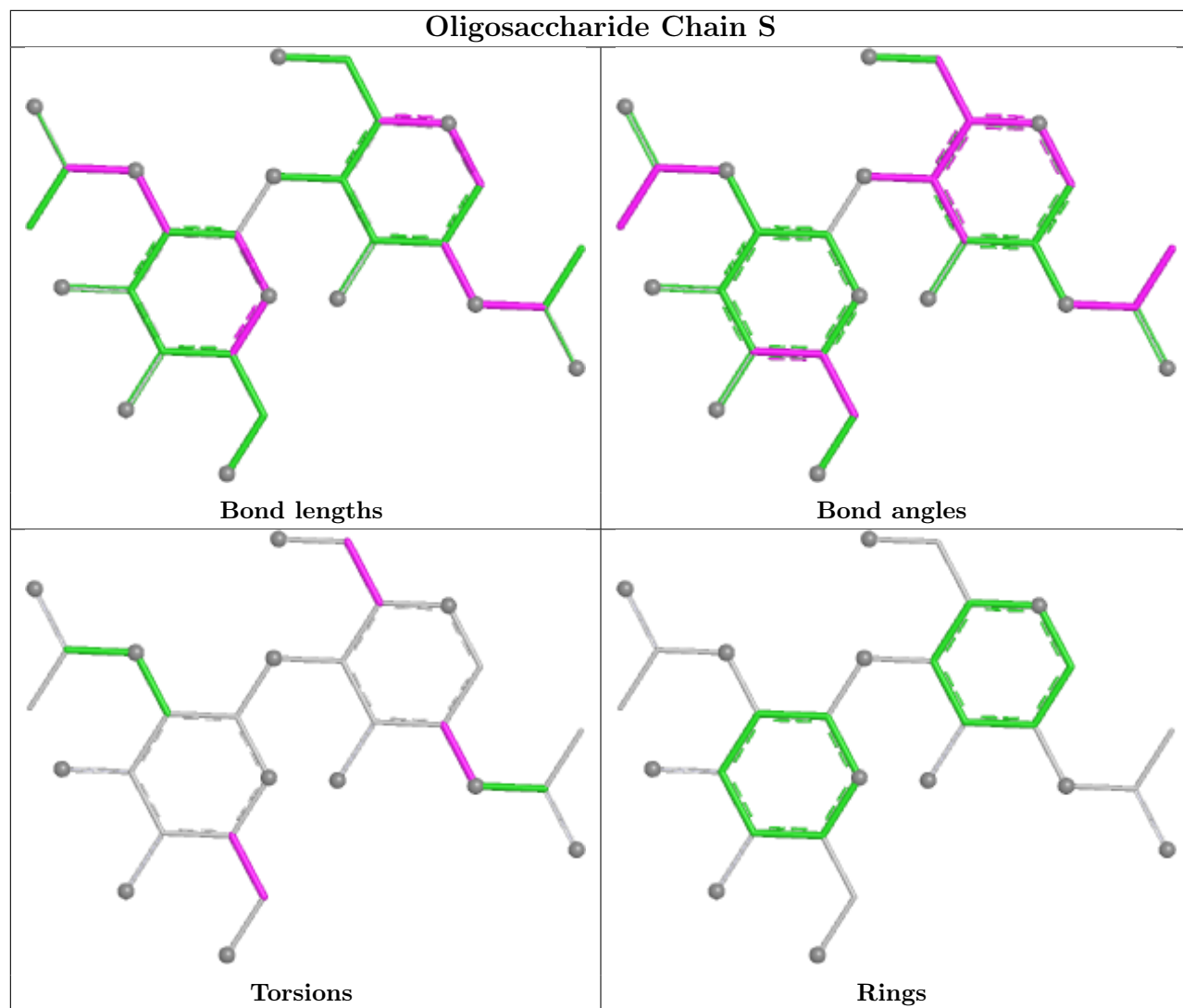


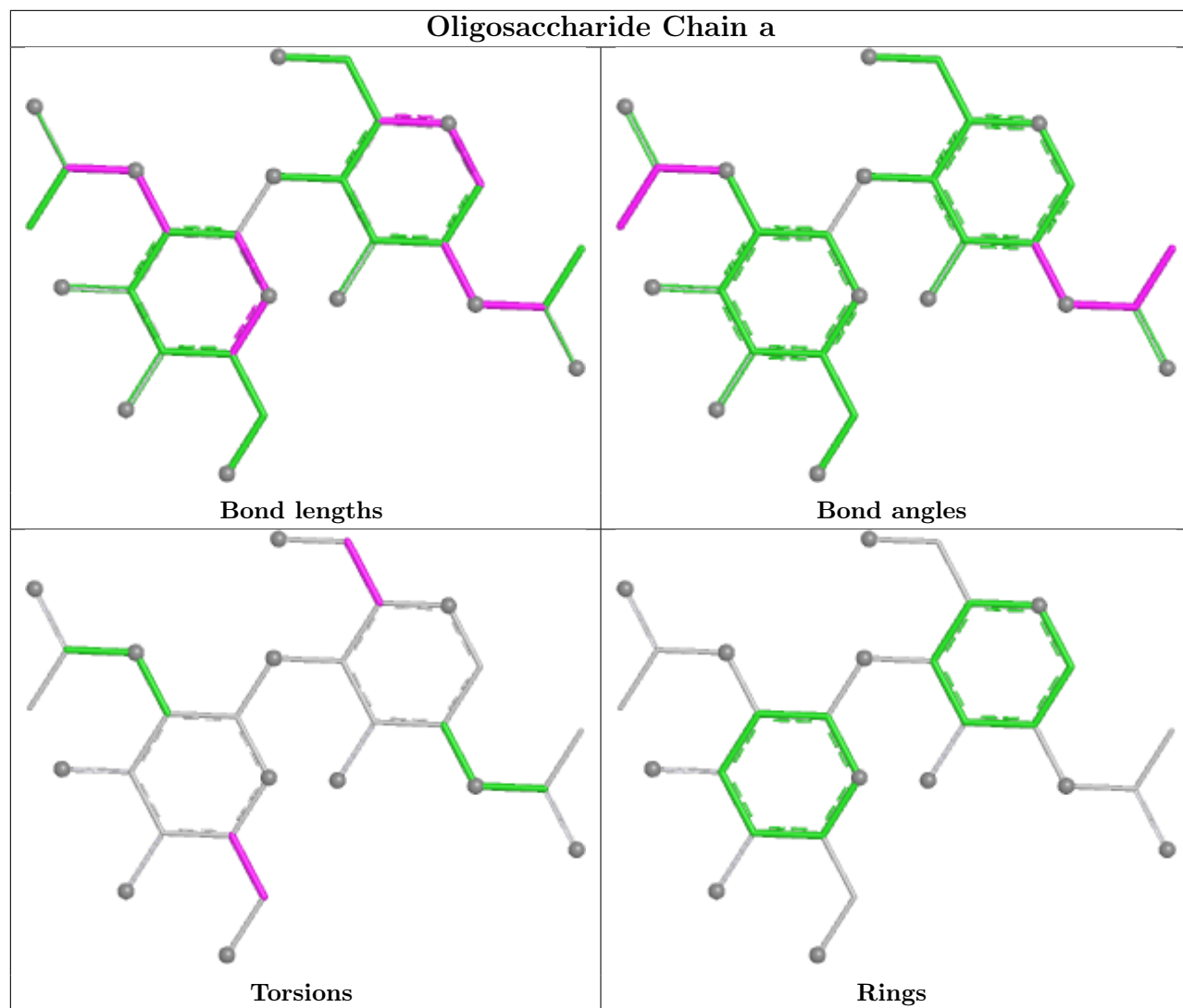


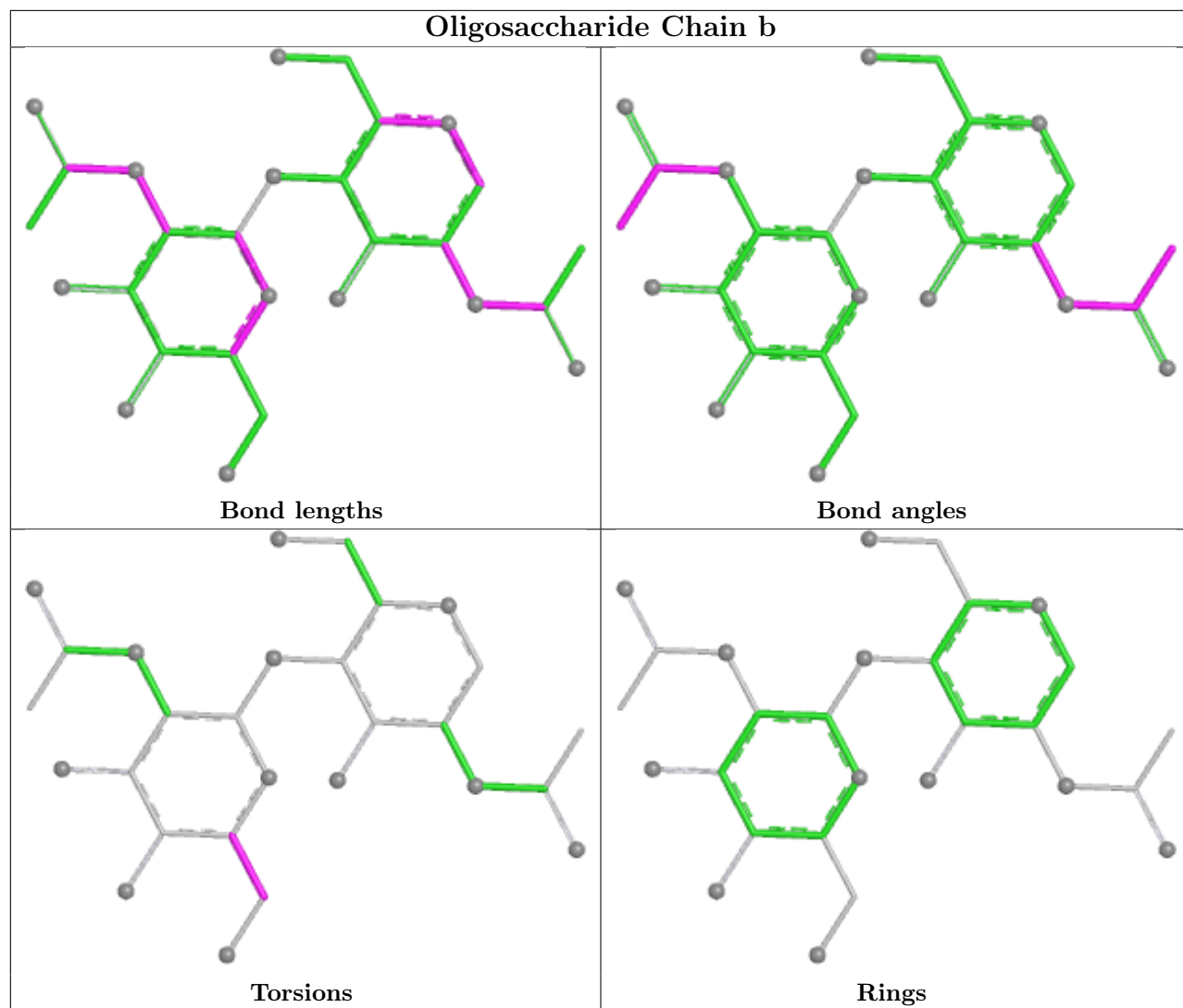


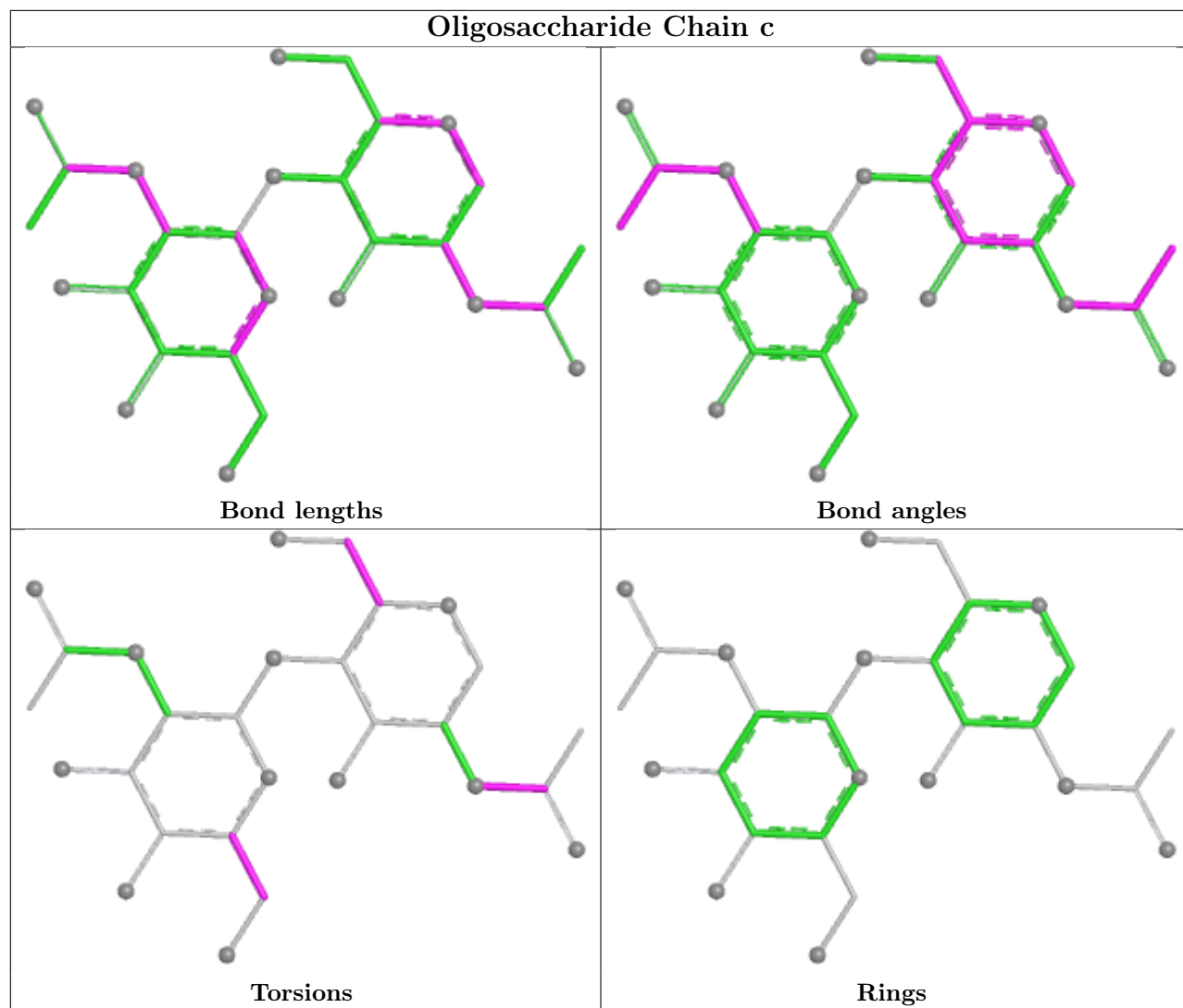


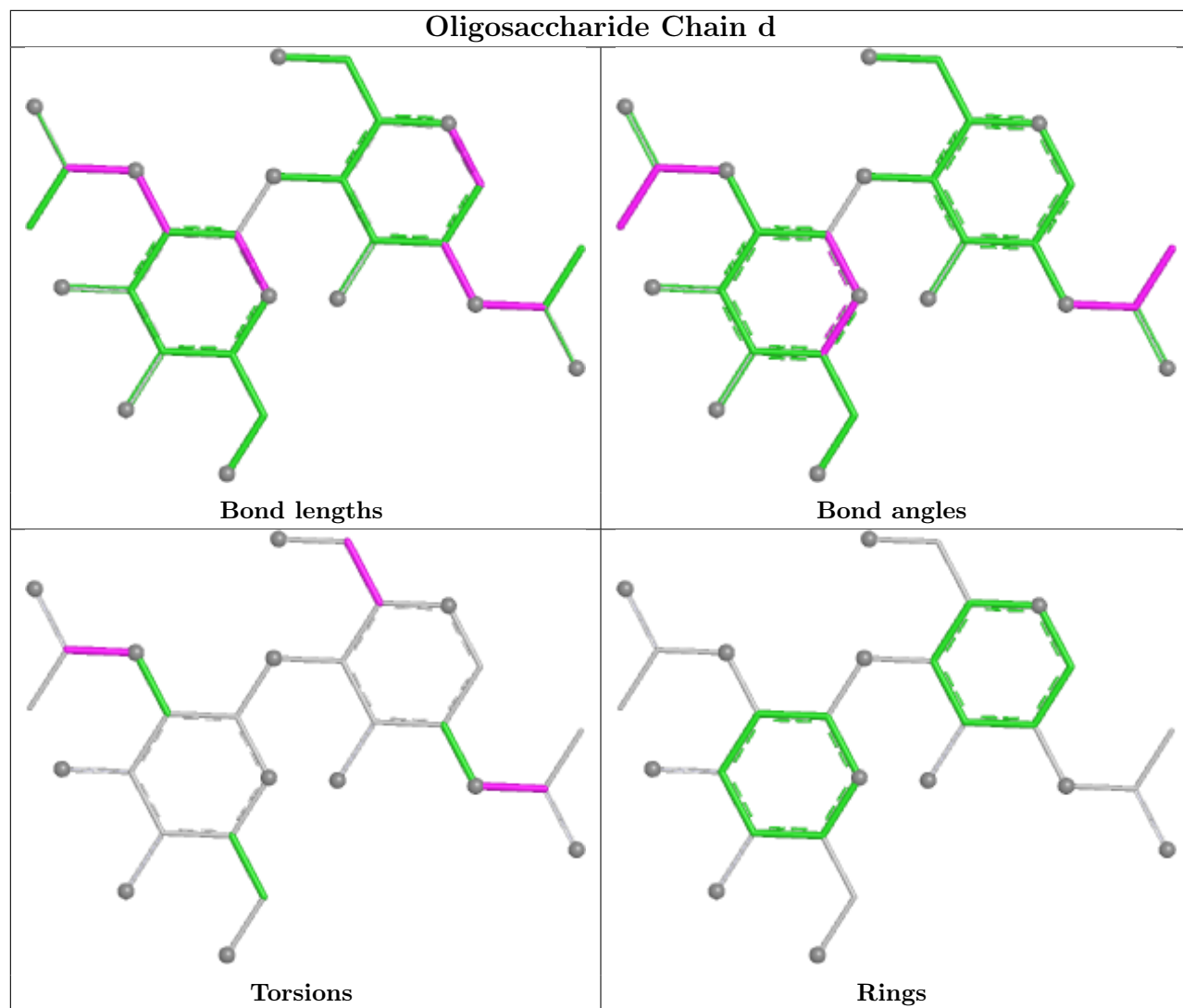


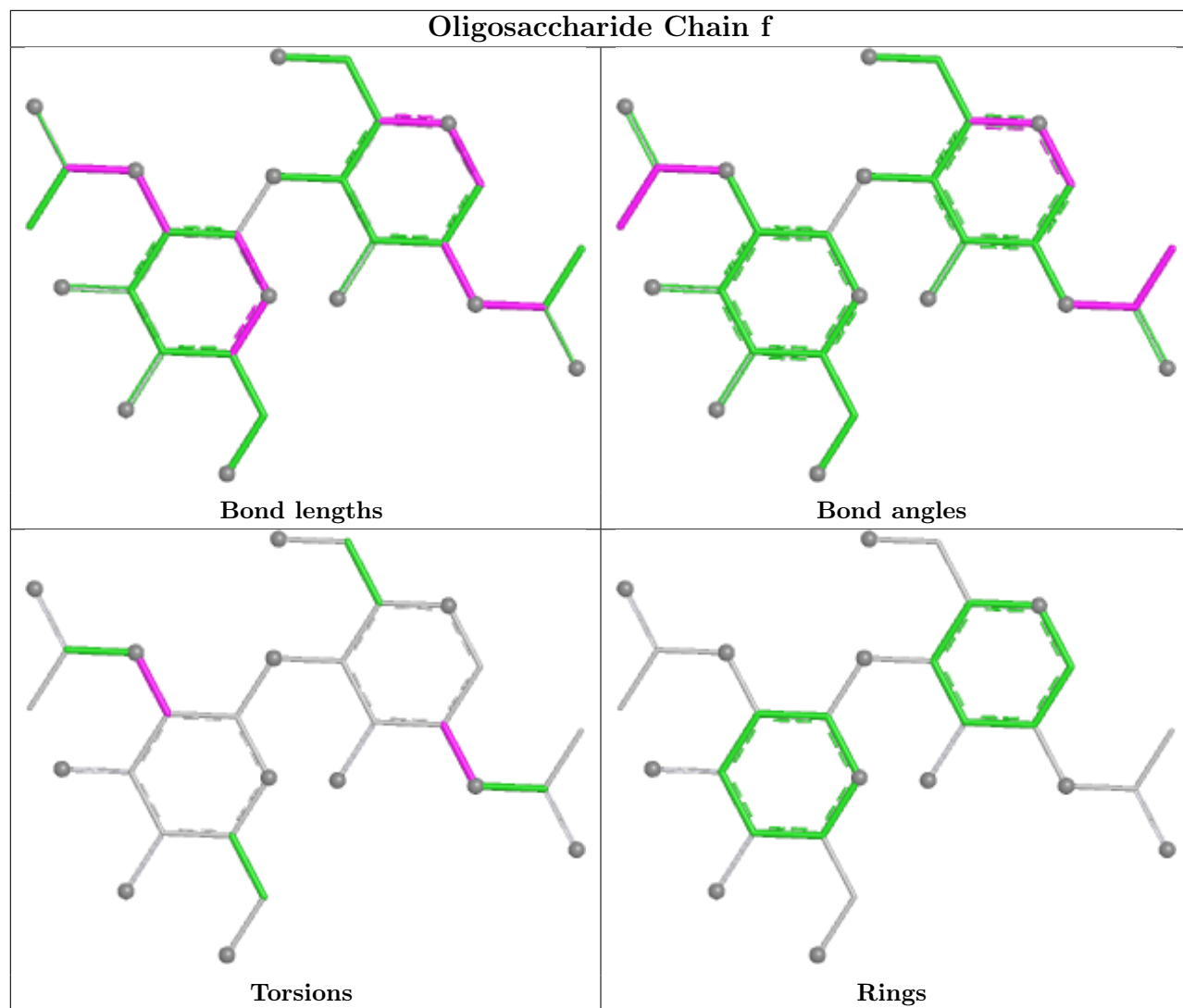


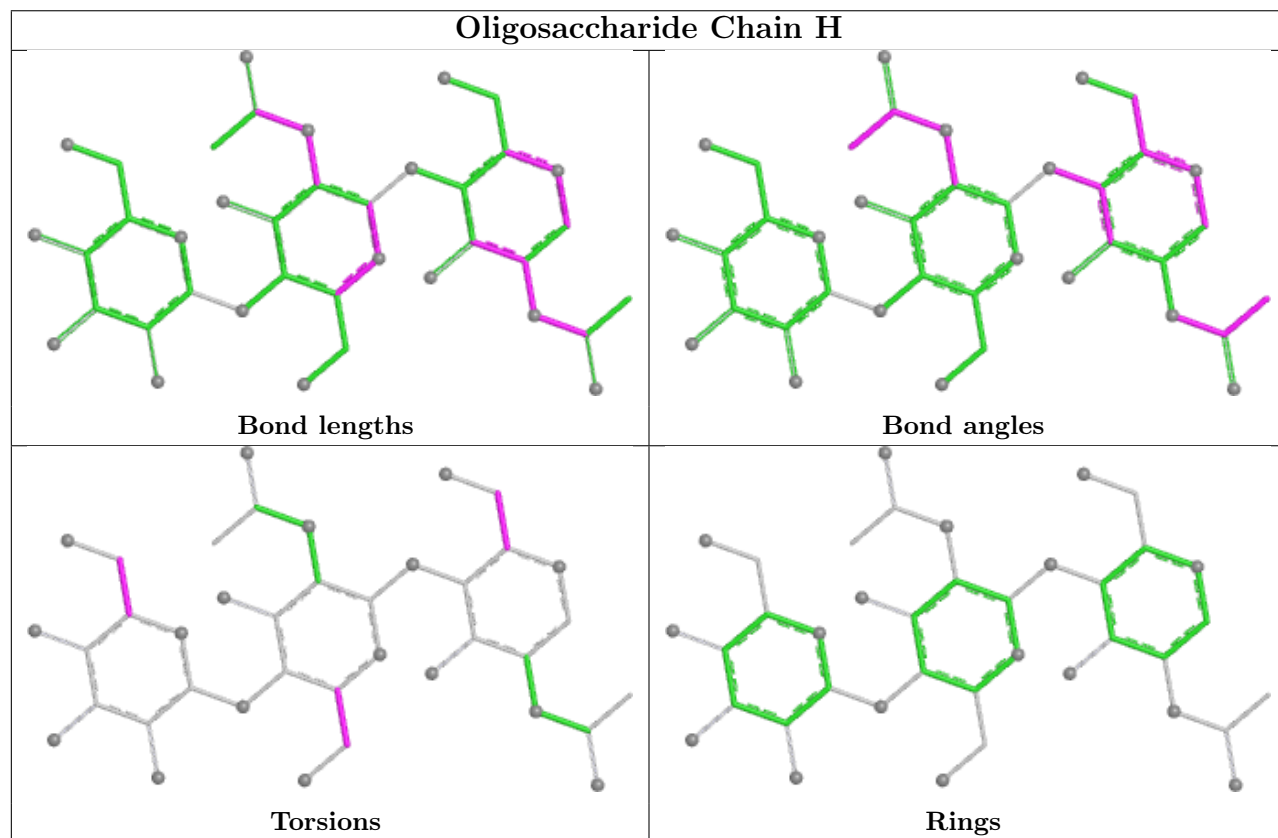
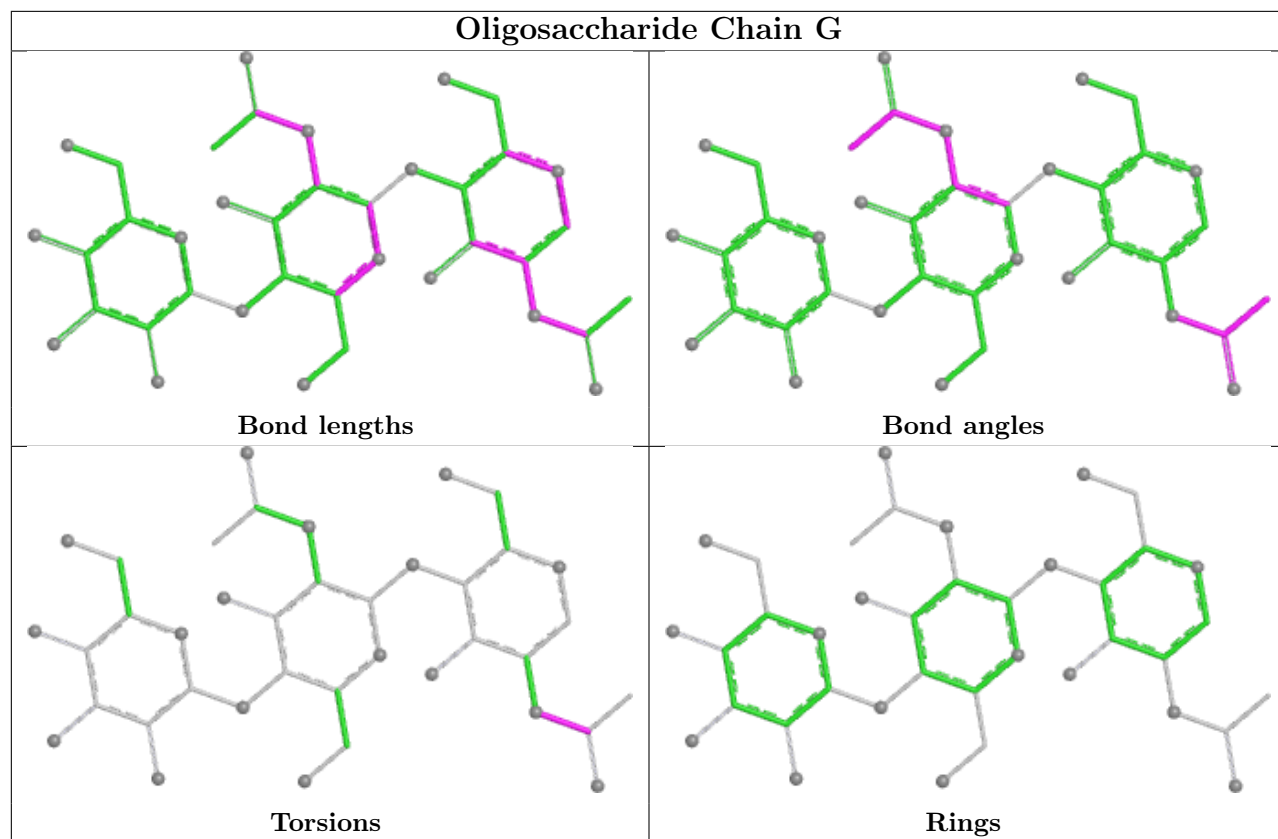


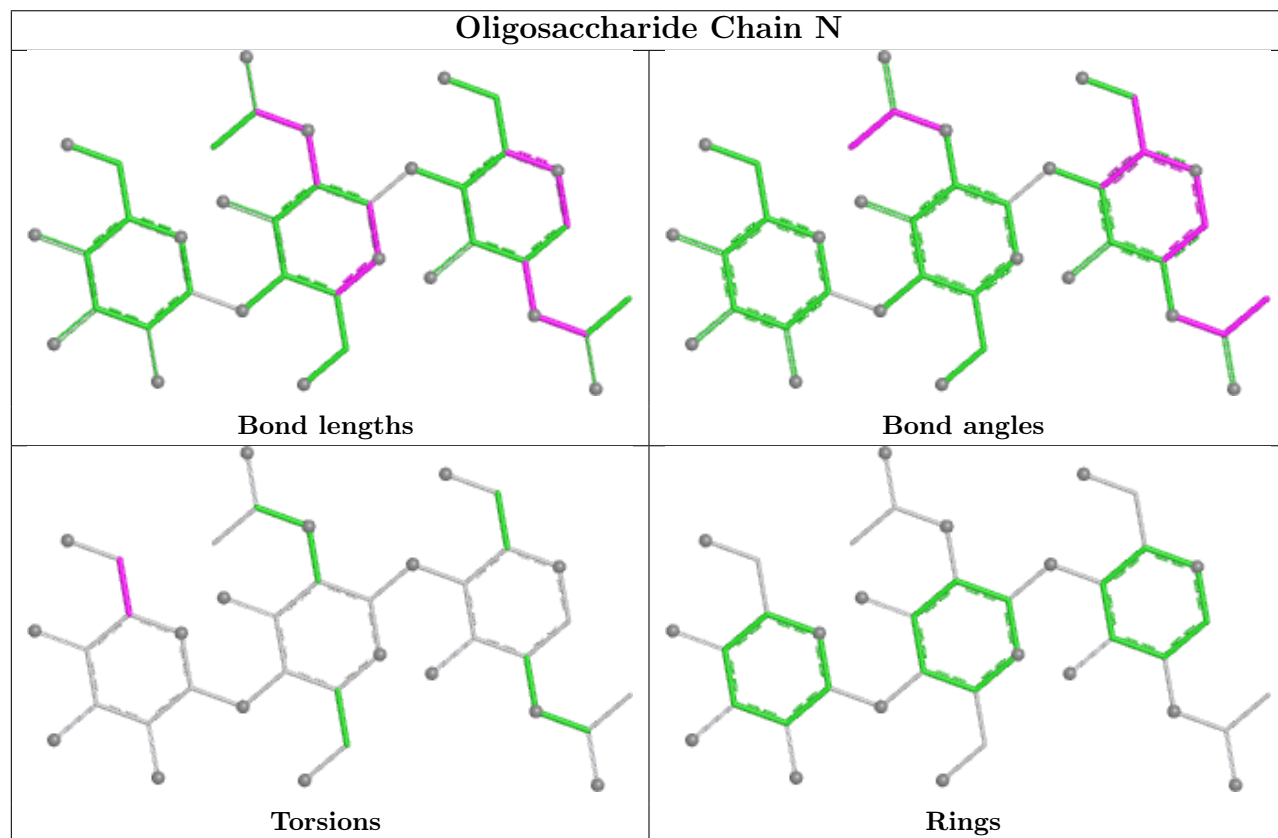
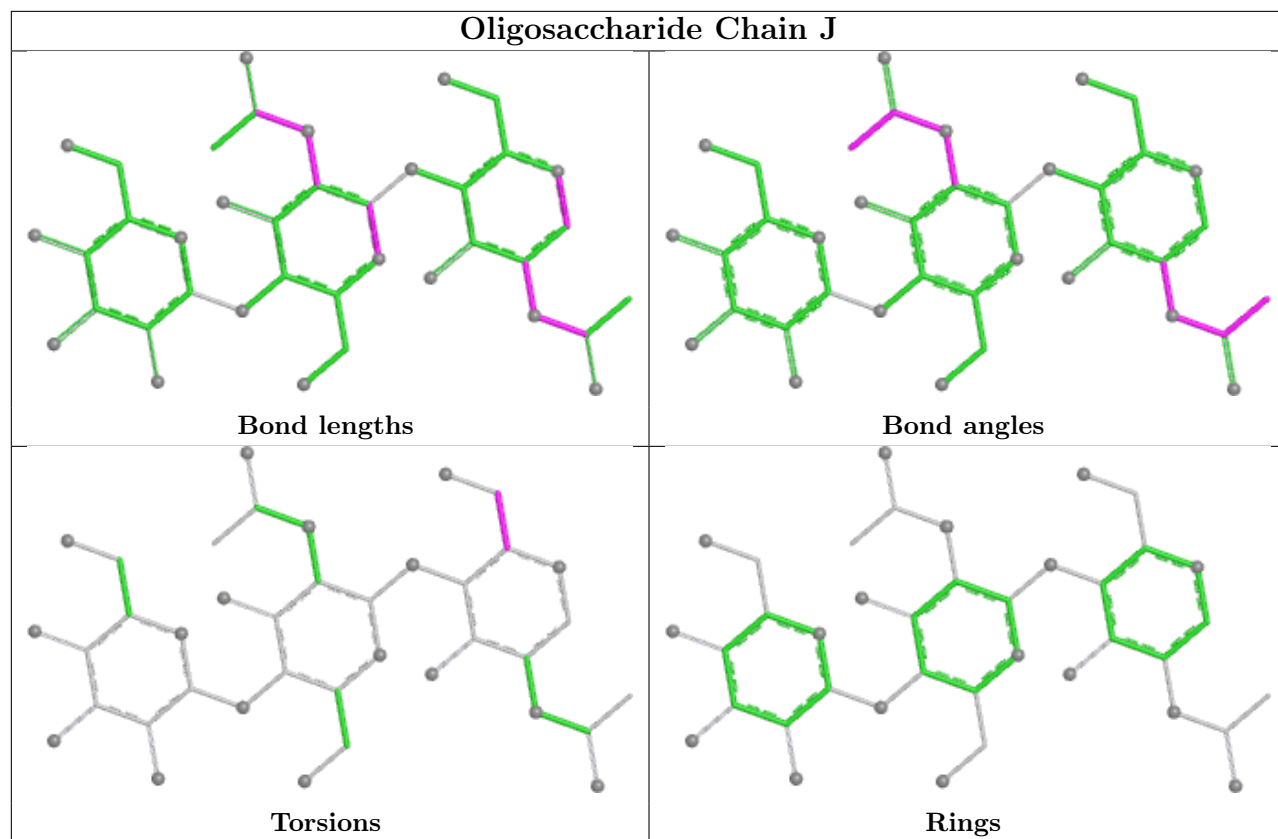


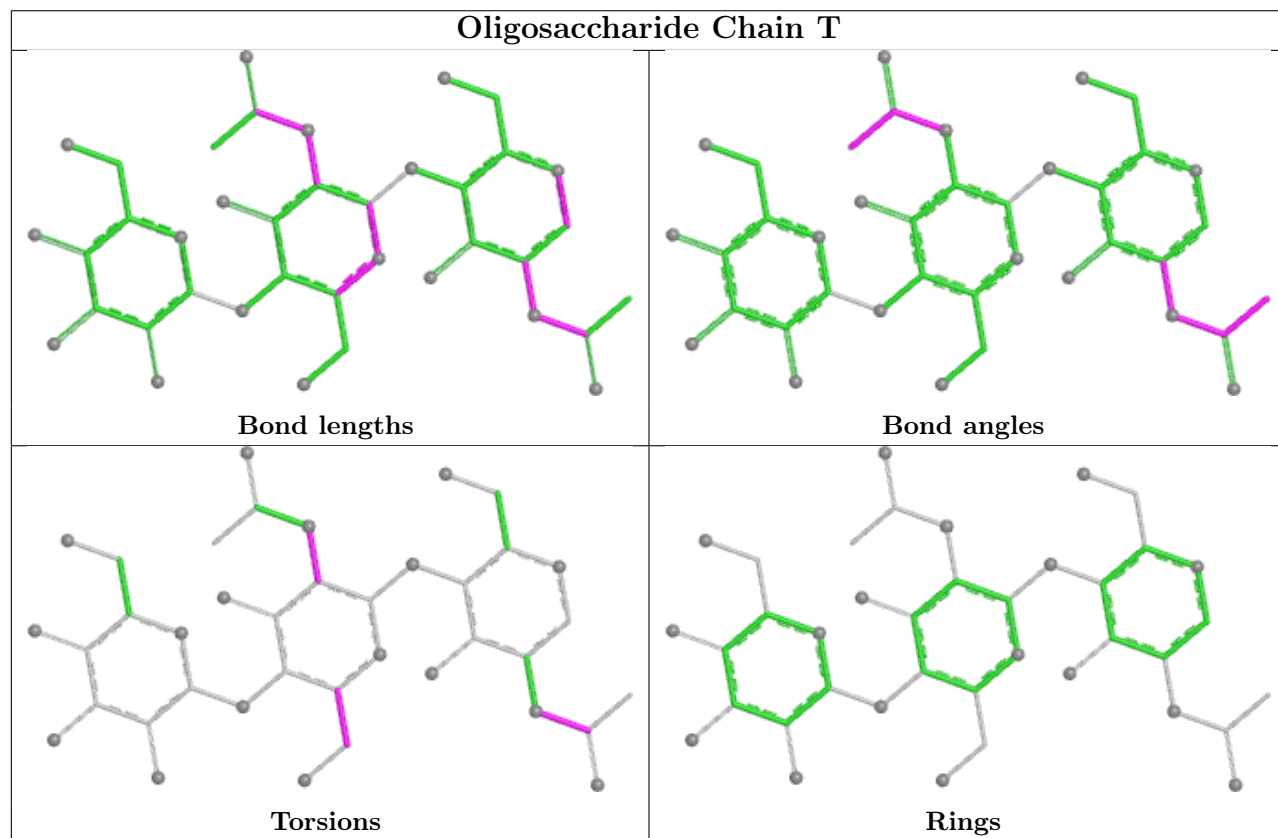
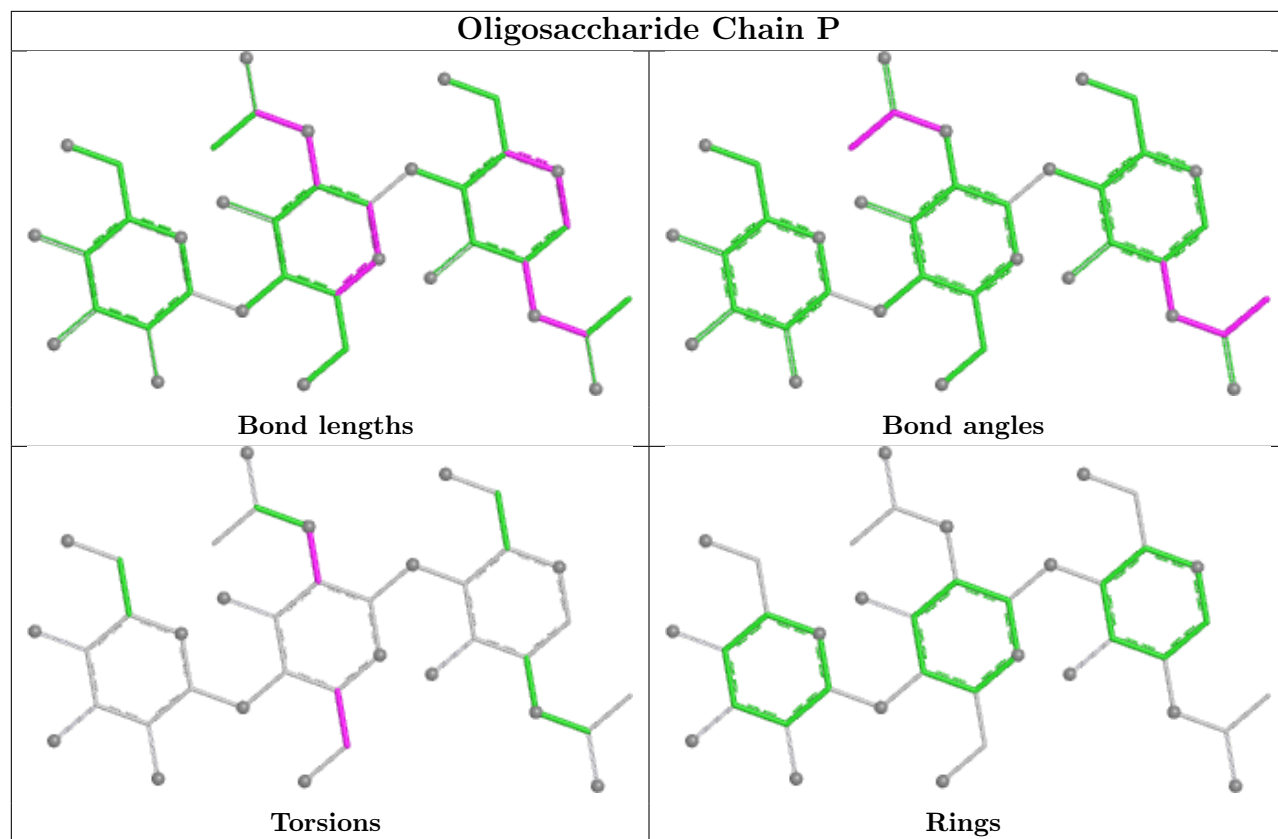


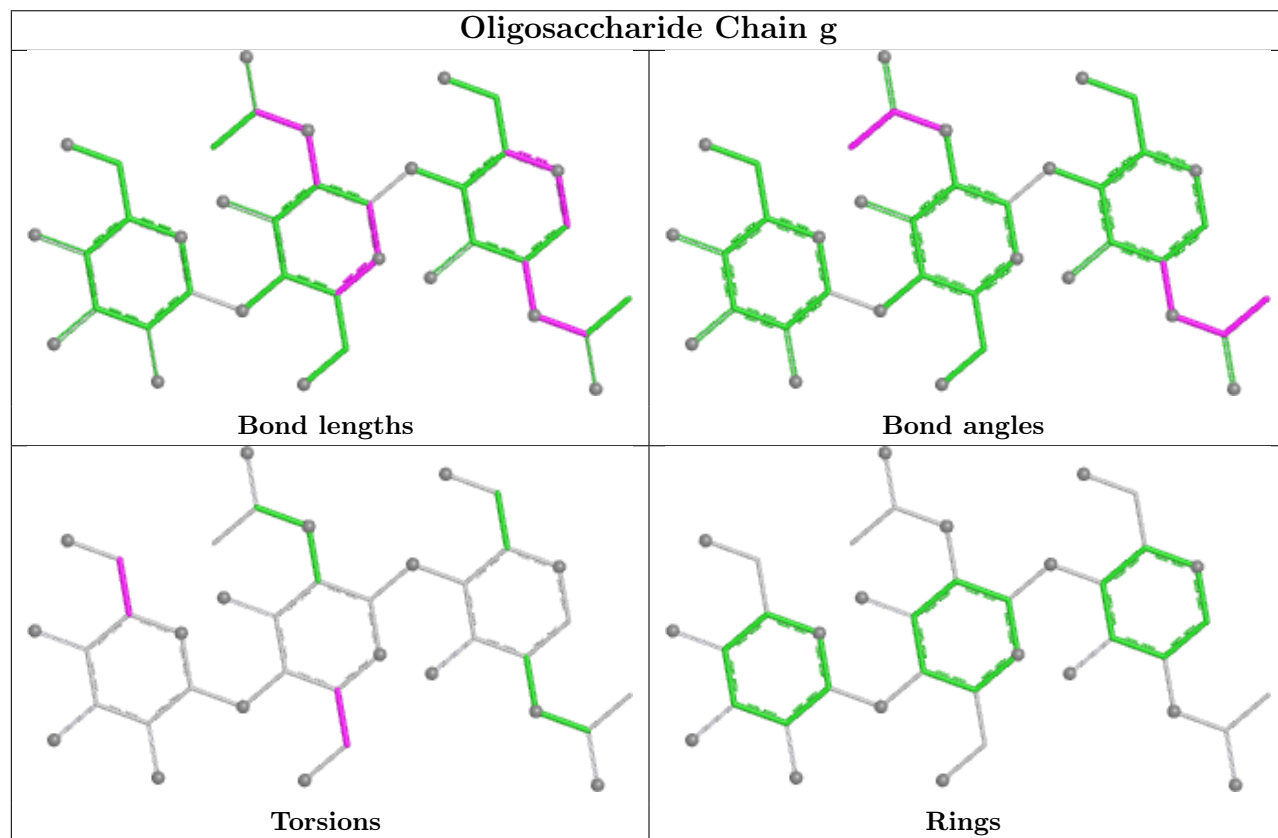
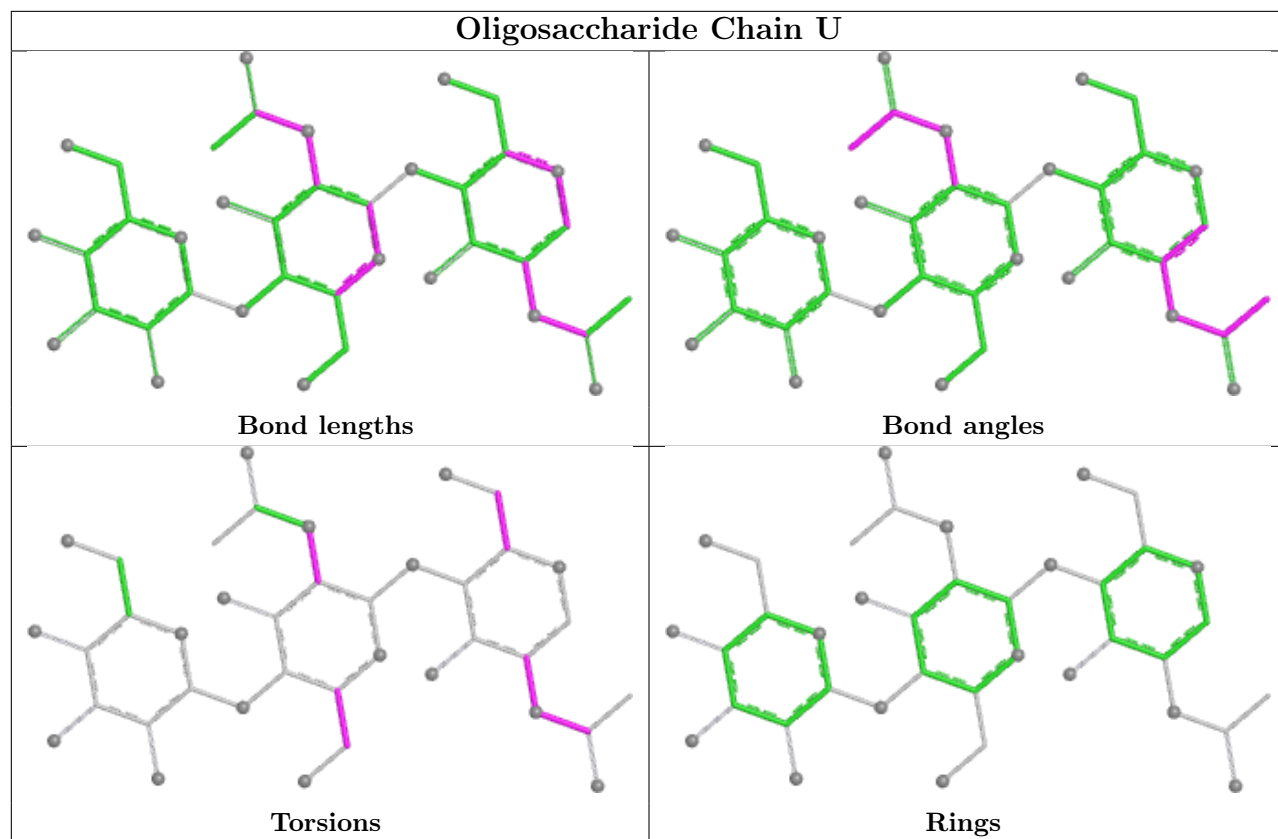


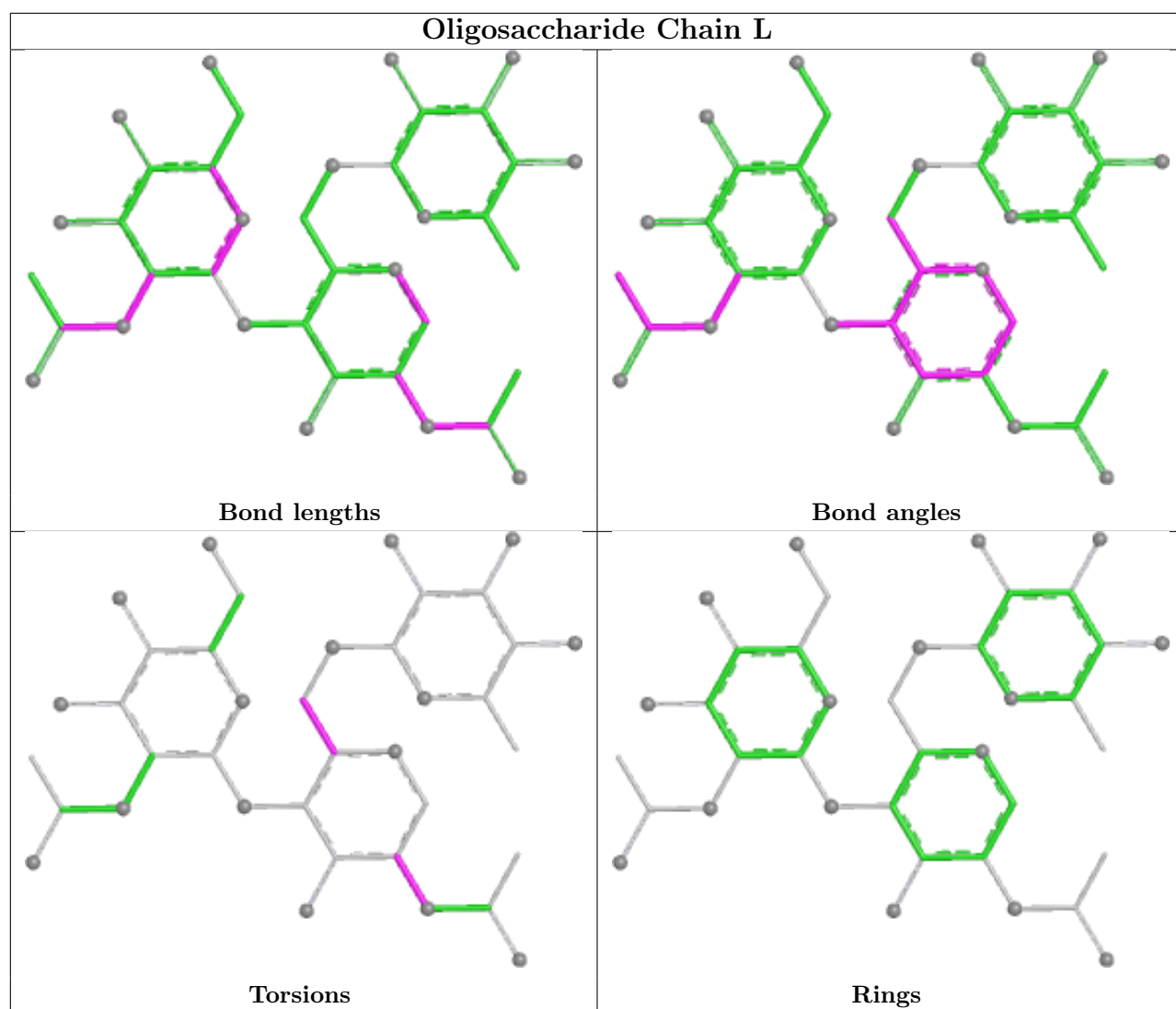


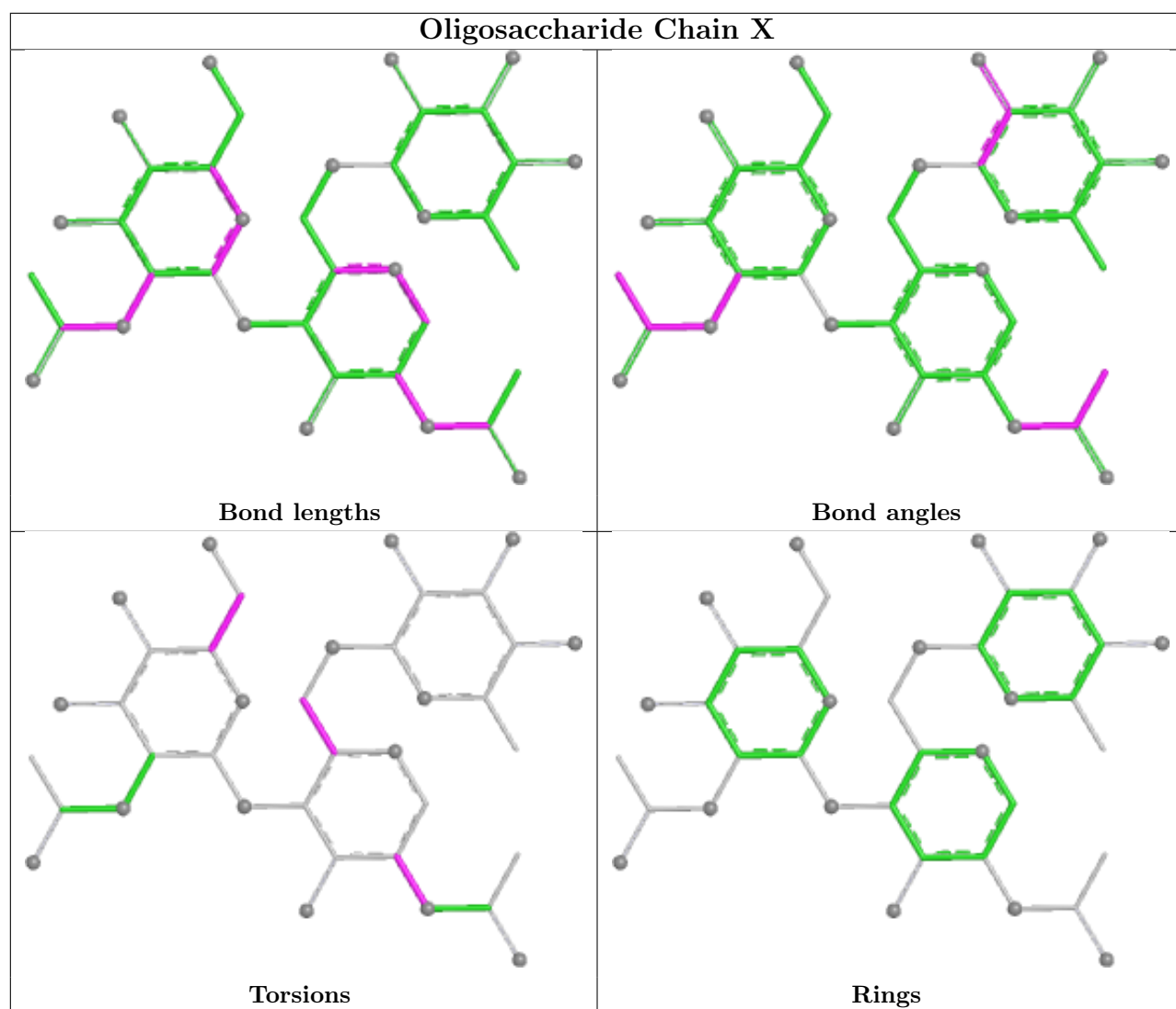


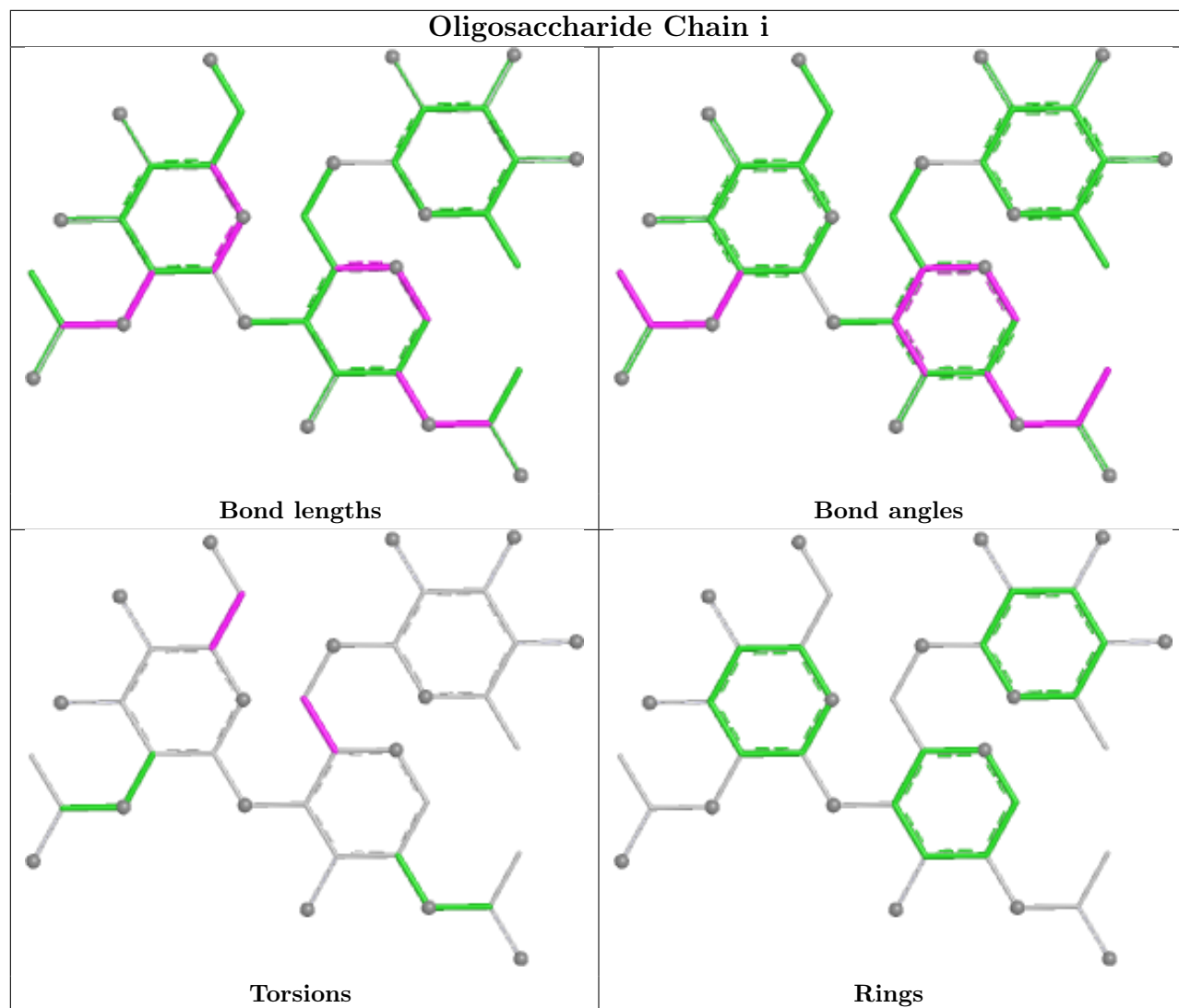


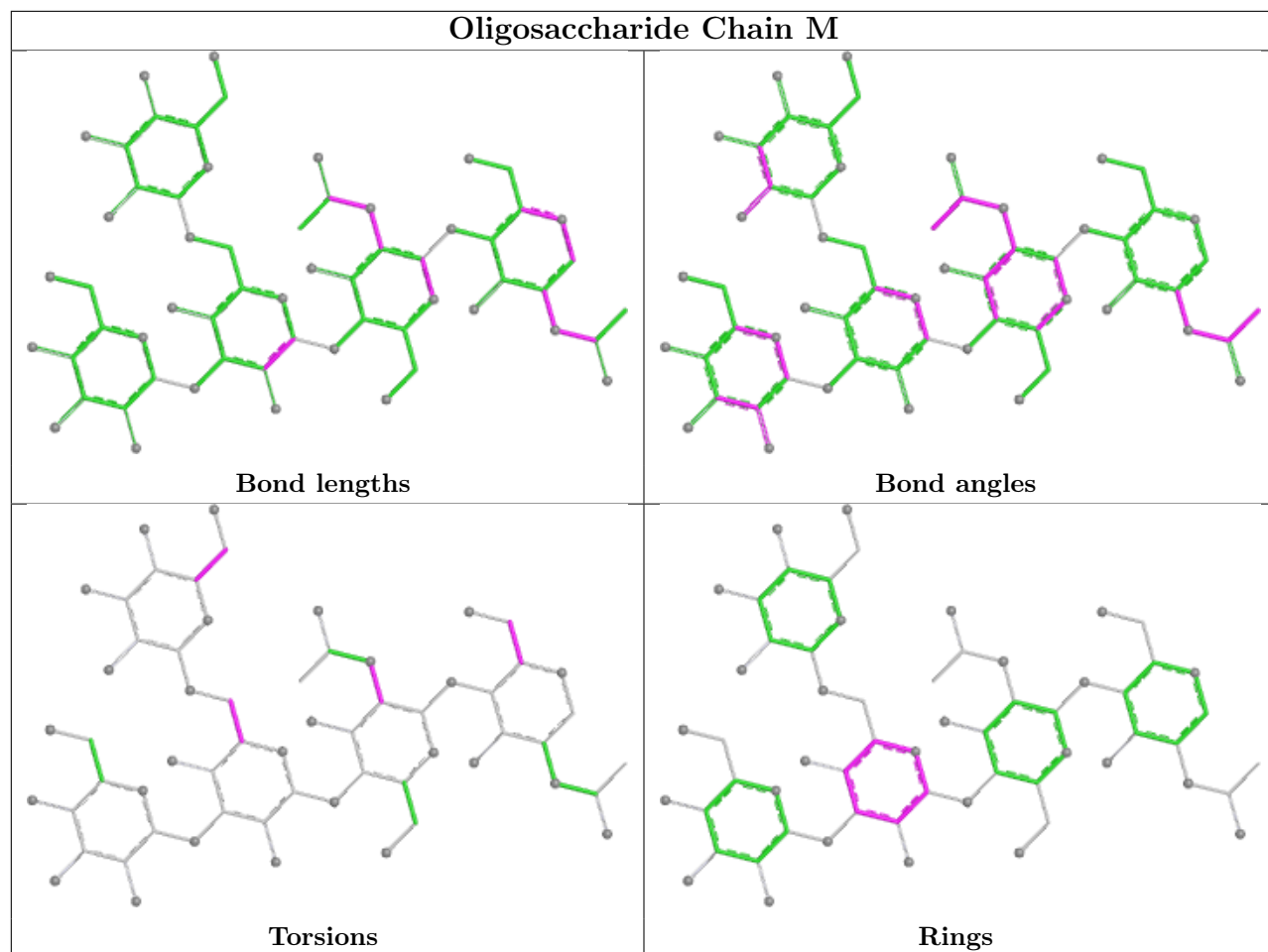


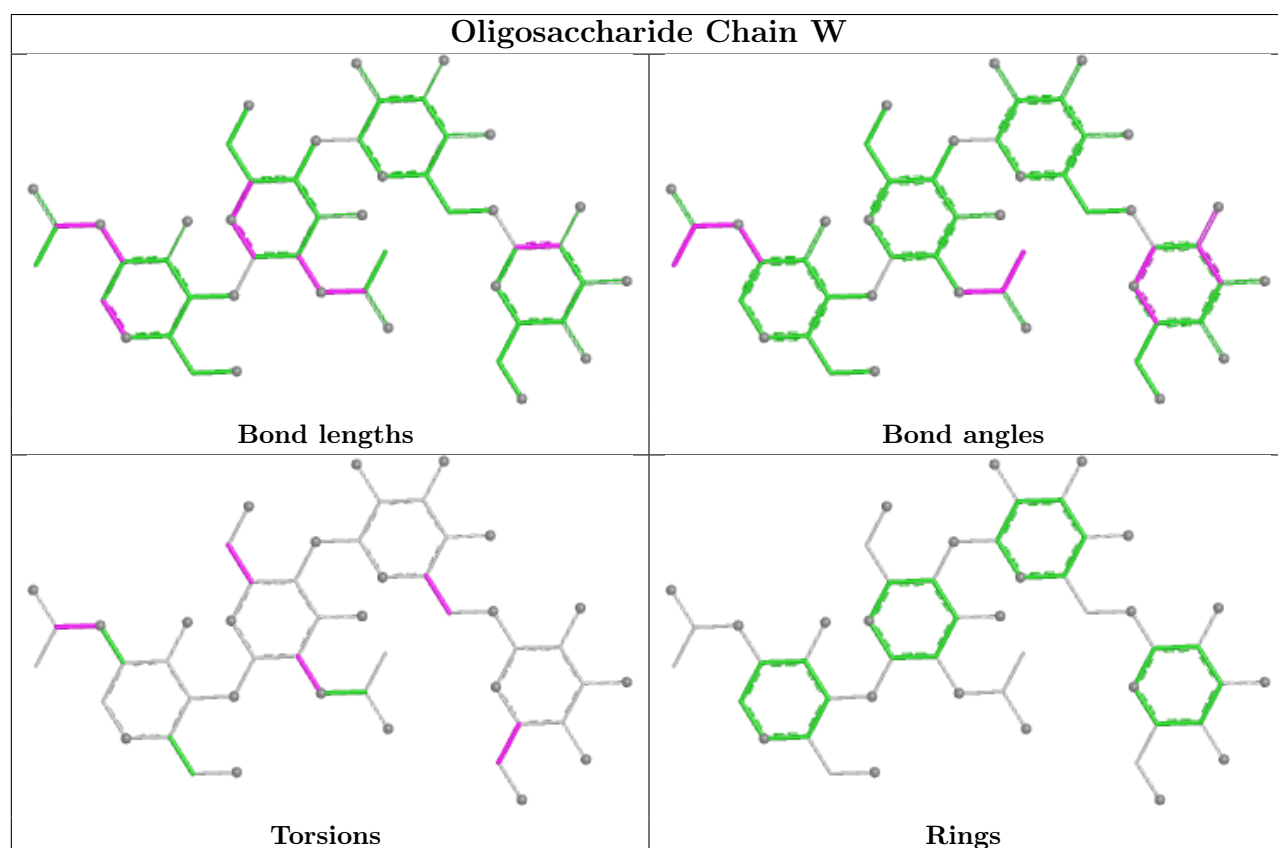
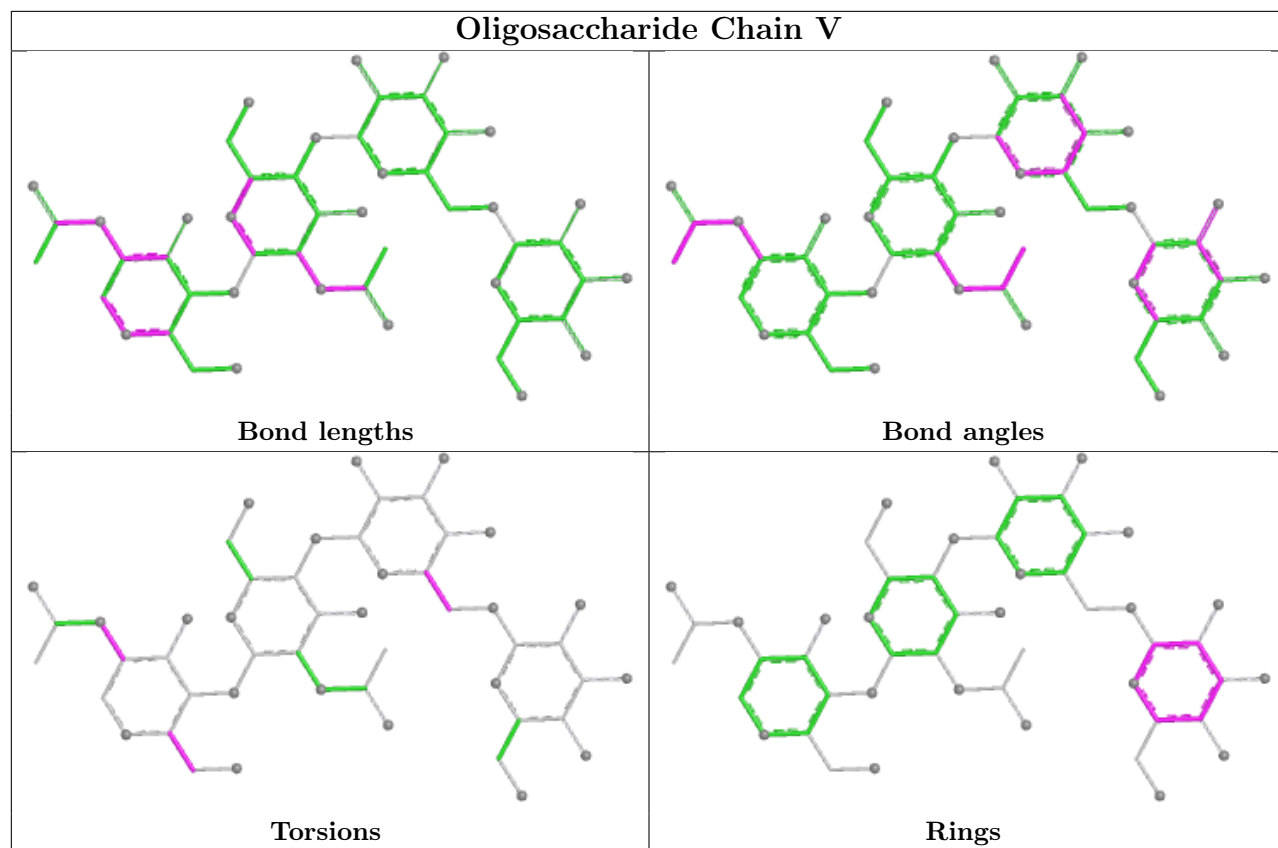


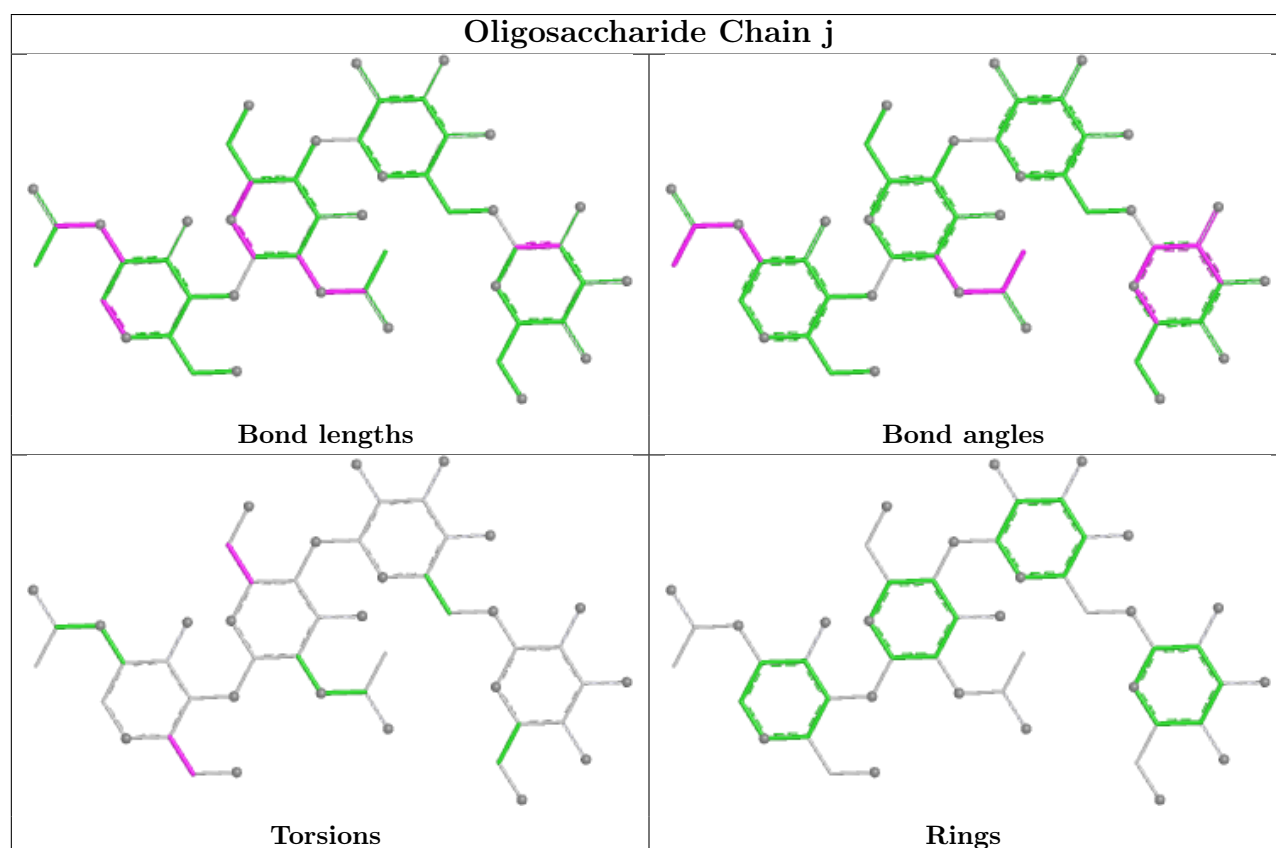
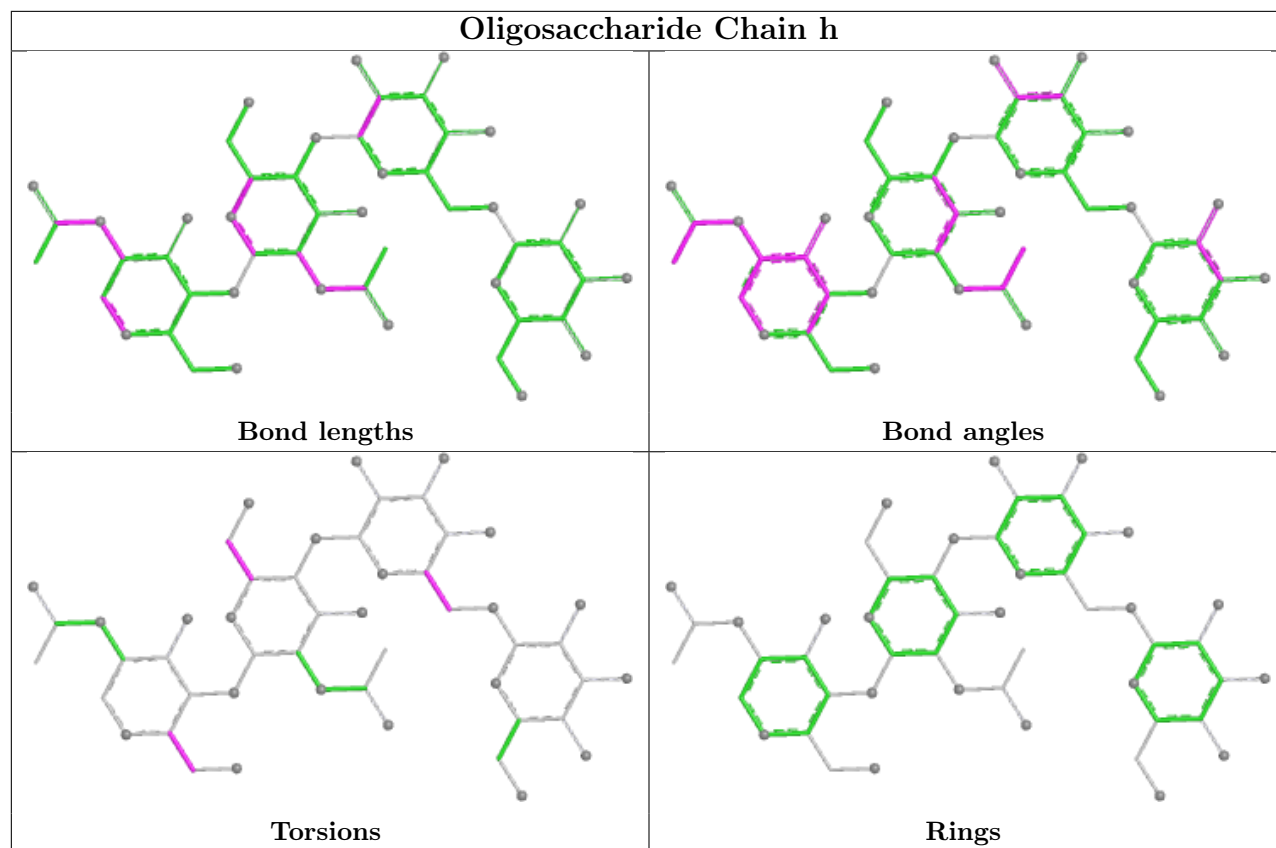


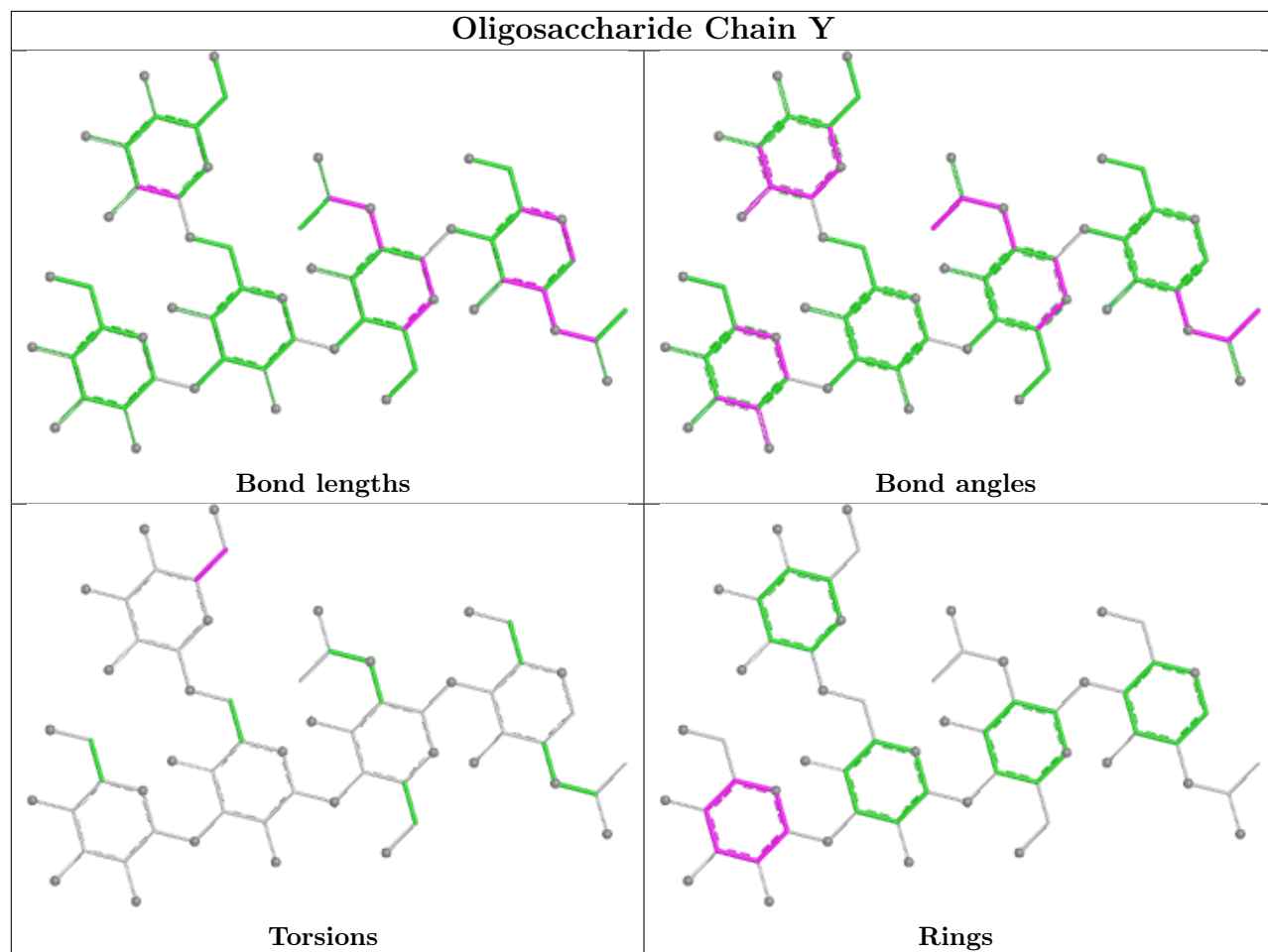


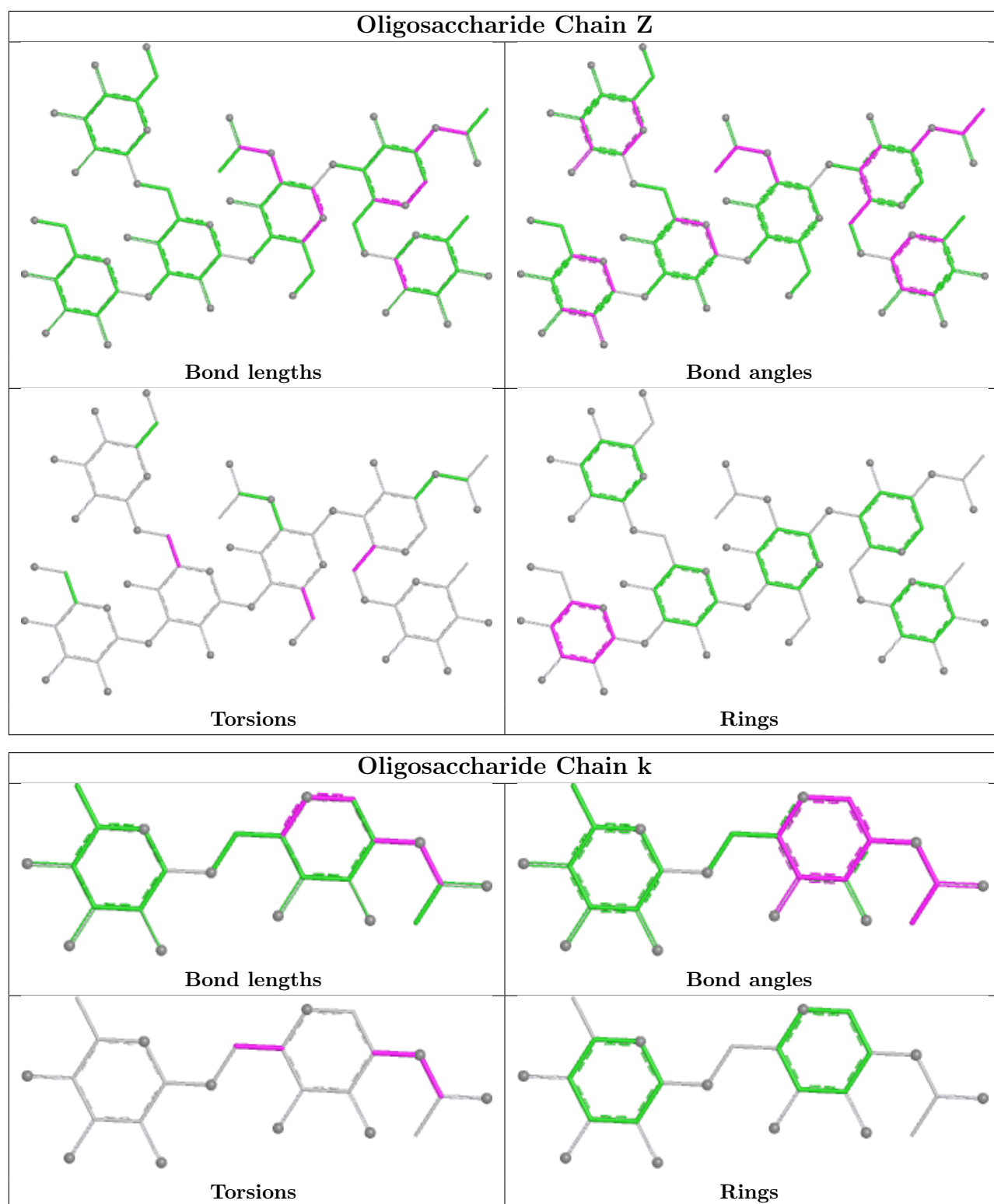


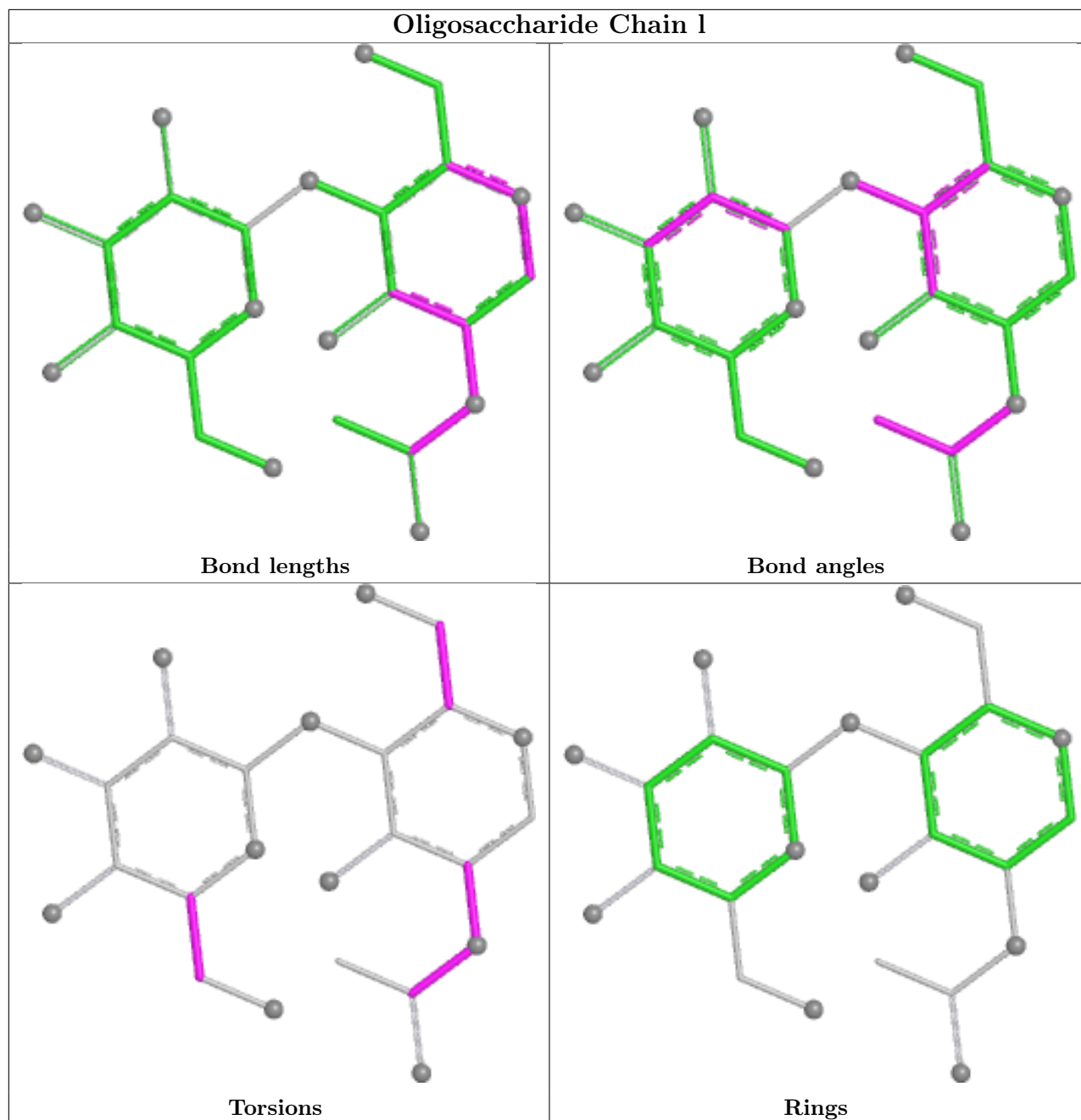












5.6 Ligand geometry [i](#)

16 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	C	1401	1	14,14,15	2.01	4 (28%)	17,19,21	1.32	1 (5%)
12	NAG	C	1402	1	14,14,15	2.02	4 (28%)	17,19,21	1.13	2 (11%)
12	NAG	C	1404	1	14,14,15	2.05	4 (28%)	17,19,21	1.20	2 (11%)
12	NAG	B	1405	1	14,14,15	1.94	3 (21%)	17,19,21	1.65	3 (17%)
12	NAG	A	1403	1	14,14,15	1.99	4 (28%)	17,19,21	1.14	2 (11%)
12	NAG	A	1405	1	14,14,15	2.05	4 (28%)	17,19,21	1.19	1 (5%)
12	NAG	B	1404	1	14,14,15	1.98	4 (28%)	17,19,21	1.21	2 (11%)
12	NAG	A	1406	-	14,14,15	1.99	4 (28%)	17,19,21	0.91	1 (5%)
12	NAG	B	1401	1	14,14,15	2.00	4 (28%)	17,19,21	1.02	1 (5%)
12	NAG	C	1403	1	14,14,15	2.02	4 (28%)	17,19,21	1.15	1 (5%)
12	NAG	C	1405	1	14,14,15	1.99	4 (28%)	17,19,21	0.97	1 (5%)
12	NAG	B	1402	1	14,14,15	1.99	4 (28%)	17,19,21	1.06	2 (11%)
12	NAG	B	1403	1	14,14,15	1.98	4 (28%)	17,19,21	1.09	2 (11%)
12	NAG	A	1402	1	14,14,15	2.01	4 (28%)	17,19,21	1.15	2 (11%)
12	NAG	A	1401	1	14,14,15	1.98	4 (28%)	17,19,21	1.13	2 (11%)
12	NAG	A	1404	1	14,14,15	2.07	4 (28%)	17,19,21	2.57	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	NAG	C	1401	1	-	3/6/23/26	0/1/1/1
12	NAG	C	1402	1	-	0/6/23/26	0/1/1/1
12	NAG	C	1404	1	-	0/6/23/26	0/1/1/1
12	NAG	B	1405	1	-	4/6/23/26	0/1/1/1
12	NAG	A	1403	1	-	2/6/23/26	0/1/1/1
12	NAG	A	1405	1	-	2/6/23/26	0/1/1/1
12	NAG	B	1404	1	-	0/6/23/26	0/1/1/1
12	NAG	A	1406	-	-	4/6/23/26	0/1/1/1
12	NAG	B	1401	1	-	2/6/23/26	0/1/1/1
12	NAG	C	1403	1	-	4/6/23/26	0/1/1/1
12	NAG	C	1405	1	-	0/6/23/26	0/1/1/1
12	NAG	B	1402	1	-	0/6/23/26	0/1/1/1
12	NAG	B	1403	1	-	1/6/23/26	0/1/1/1
12	NAG	A	1402	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	NAG	A	1401	1	-	0/6/23/26	0/1/1/1
12	NAG	A	1404	1	1/1/6/7	5/6/23/26	0/1/1/1

The worst 5 of 63 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	C	1404	NAG	O5-C1	4.51	1.51	1.43
12	C	1402	NAG	O5-C1	4.41	1.51	1.43
12	A	1405	NAG	O5-C1	4.39	1.51	1.43
12	C	1405	NAG	O5-C1	4.37	1.51	1.43
12	A	1402	NAG	O5-C1	4.33	1.51	1.43

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	A	1404	NAG	C3-C4-C5	6.03	121.17	110.23
12	A	1404	NAG	C8-C7-N2	4.44	123.48	116.12
12	A	1404	NAG	O4-C4-C5	3.82	118.74	109.32
12	A	1404	NAG	C1-O5-C5	-3.76	107.14	112.19
12	B	1405	NAG	C4-C3-C2	3.75	116.52	111.02

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
12	A	1404	NAG	C4

5 of 29 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	A	1404	NAG	C1-C2-N2-C7
12	A	1406	NAG	C1-C2-N2-C7
12	A	1404	NAG	C4-C5-C6-O6
12	A	1403	NAG	O5-C5-C6-O6
12	B	1401	NAG	O5-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	B	1405	NAG	2	0
12	A	1406	NAG	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	A	1402	NAG	1	0
12	A	1404	NAG	5	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.

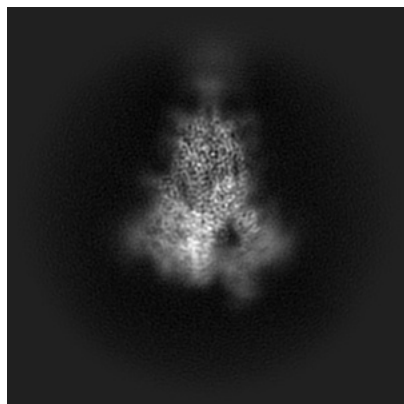
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-44447. These allow visual inspection of the internal detail of the map and identification of artifacts.

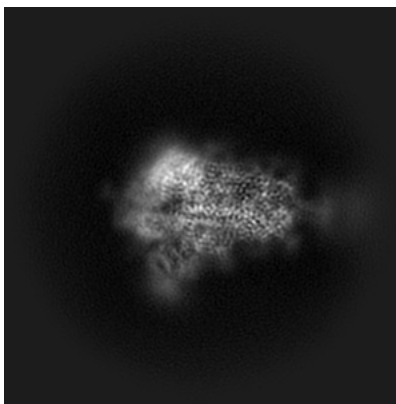
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

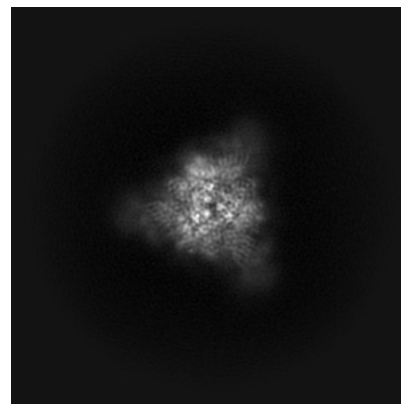
6.1.1 Primary map



X

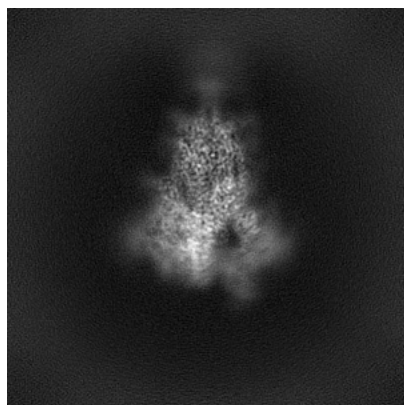


Y

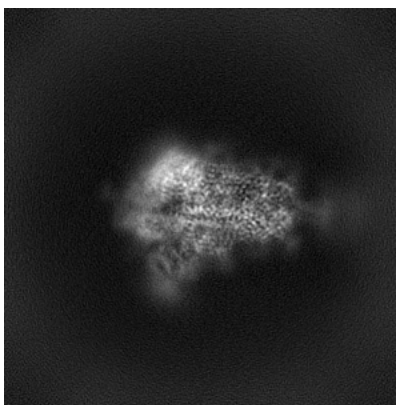


Z

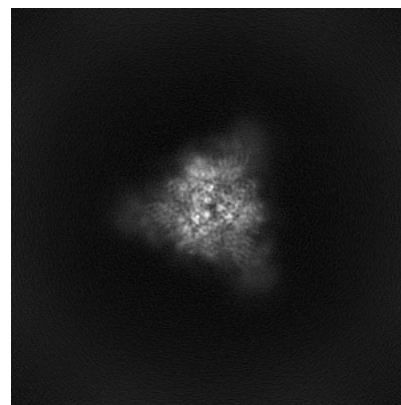
6.1.2 Raw map



X



Y

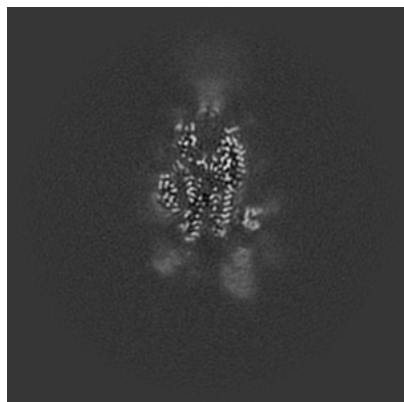


Z

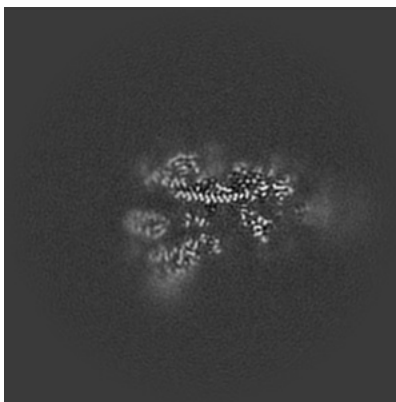
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

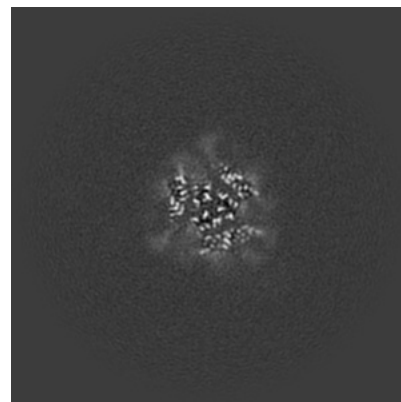
6.2.1 Primary map



X Index: 150

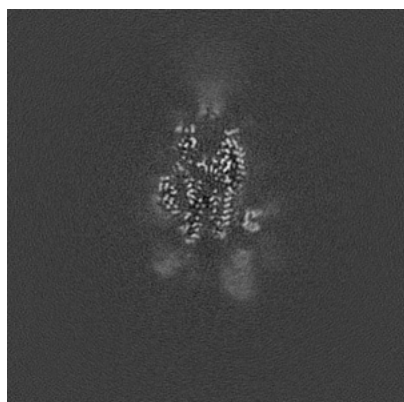


Y Index: 150

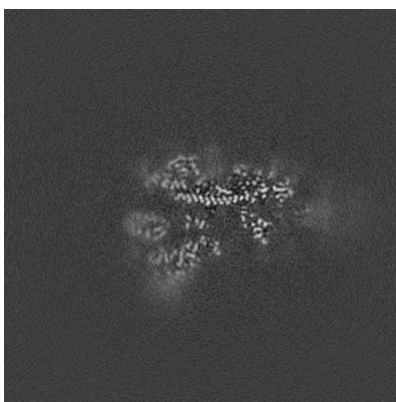


Z Index: 150

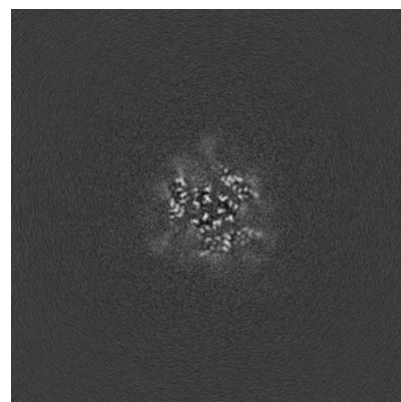
6.2.2 Raw map



X Index: 150



Y Index: 150

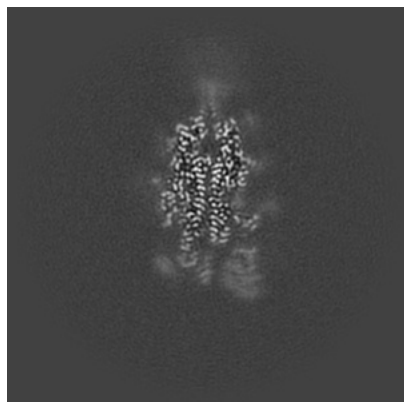


Z Index: 150

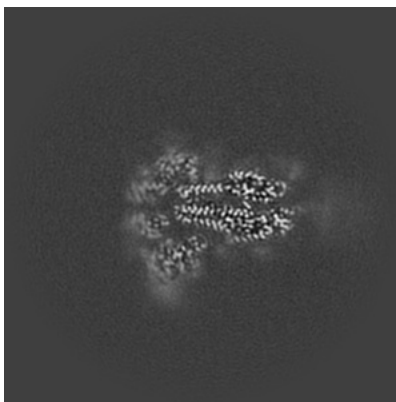
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

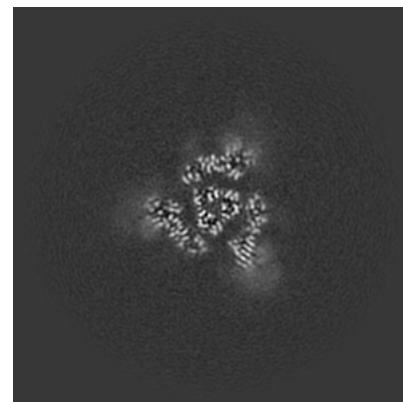
6.3.1 Primary map



X Index: 145

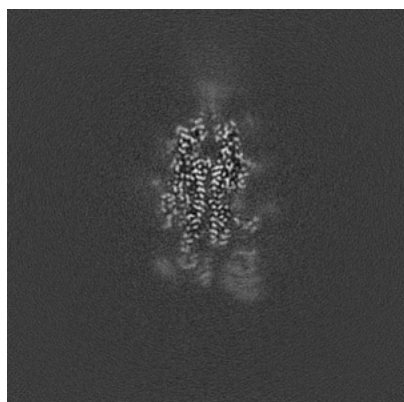


Y Index: 143

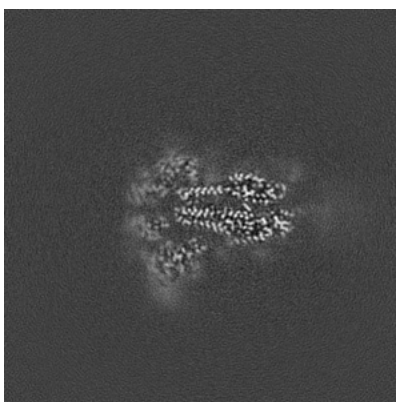


Z Index: 135

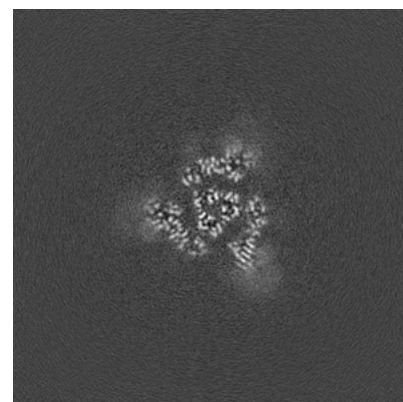
6.3.2 Raw map



X Index: 145



Y Index: 143

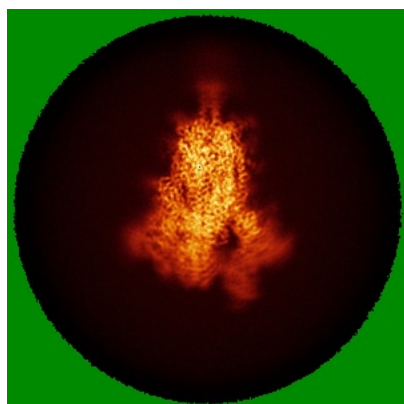


Z Index: 135

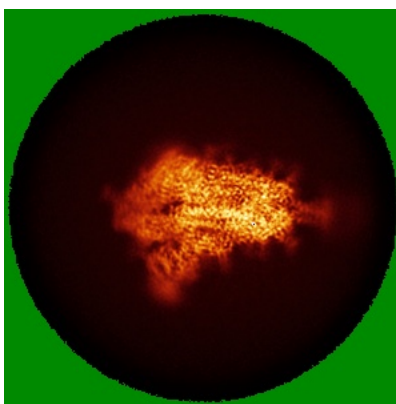
The images above show the largest variance slices of the map in three orthogonal directions.

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

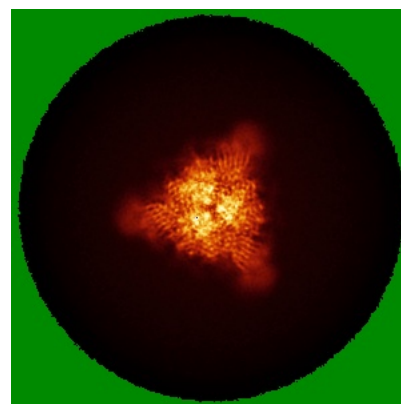
6.4.1 Primary map



X

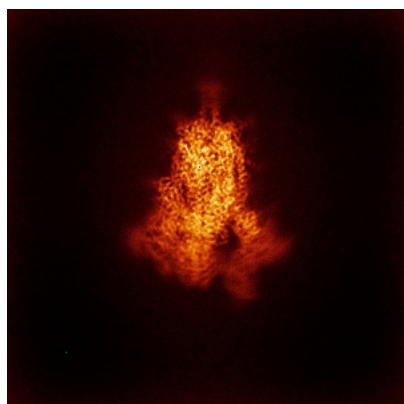


Y

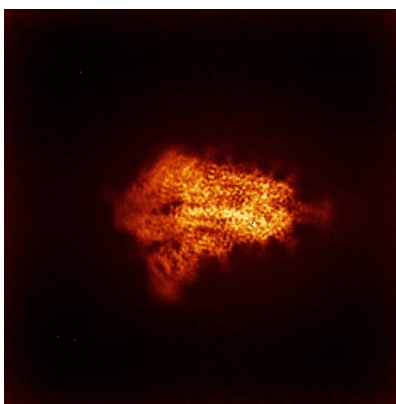


Z

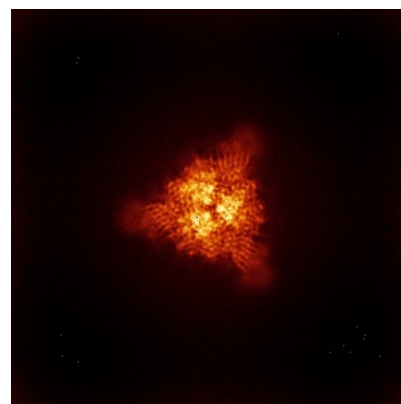
6.4.2 Raw map



X



Y

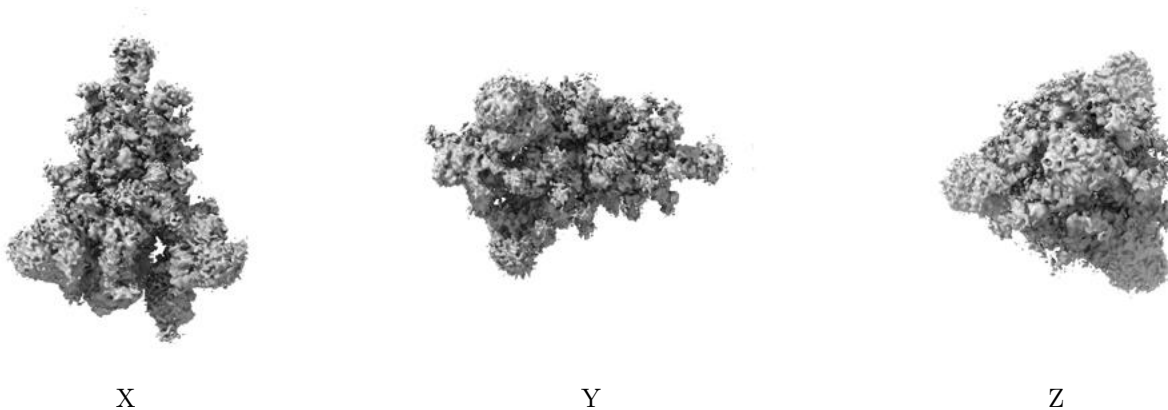


Z

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.

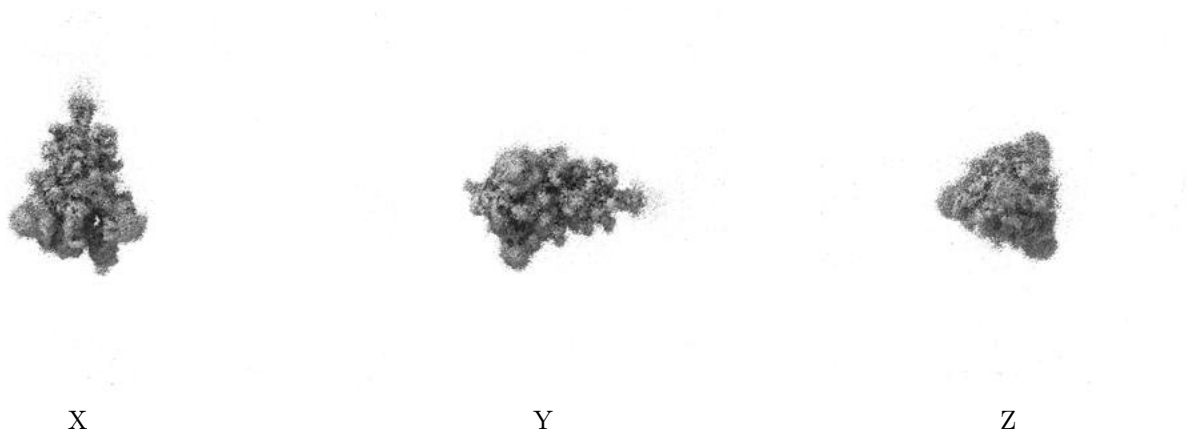
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.08. 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.

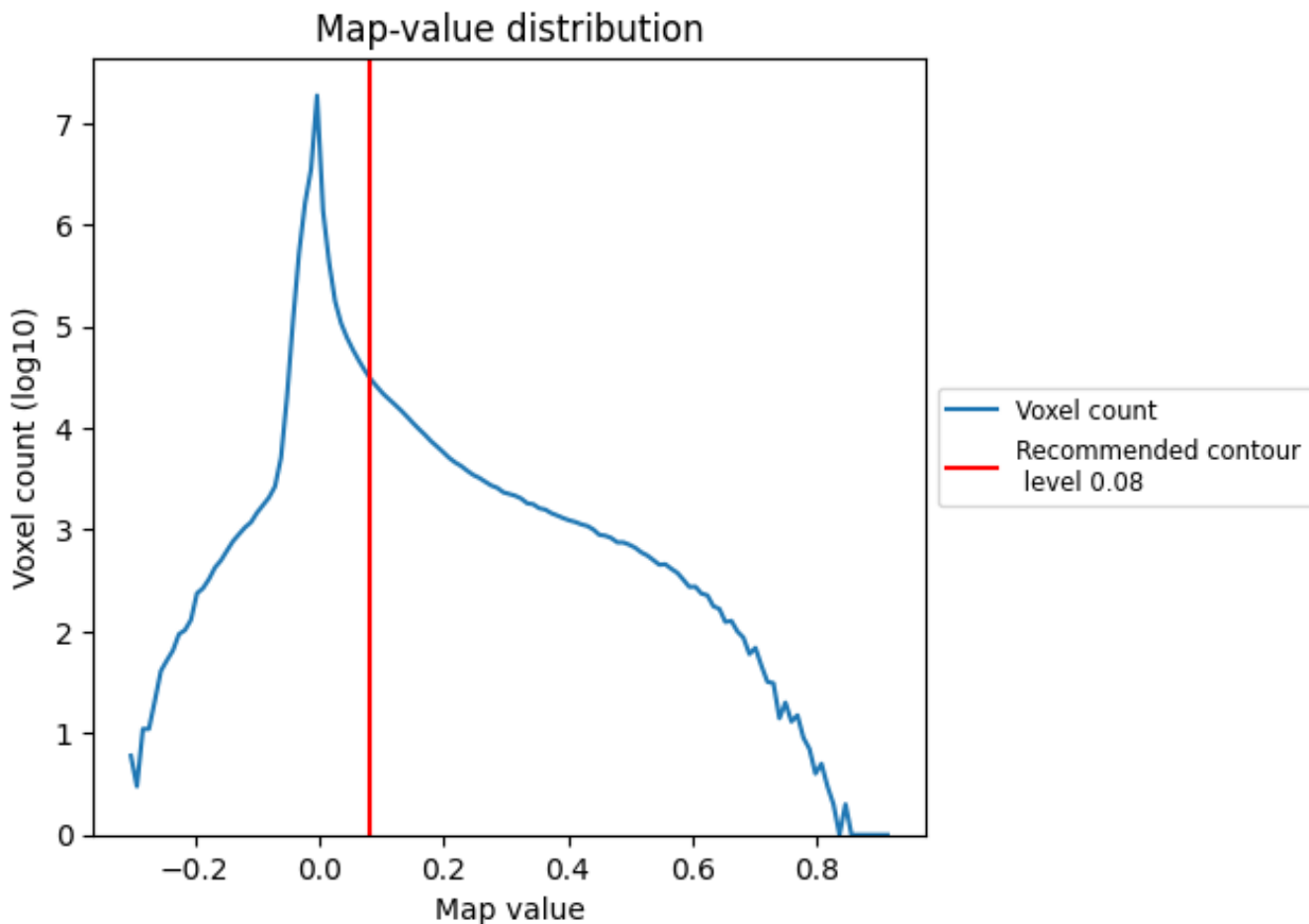
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

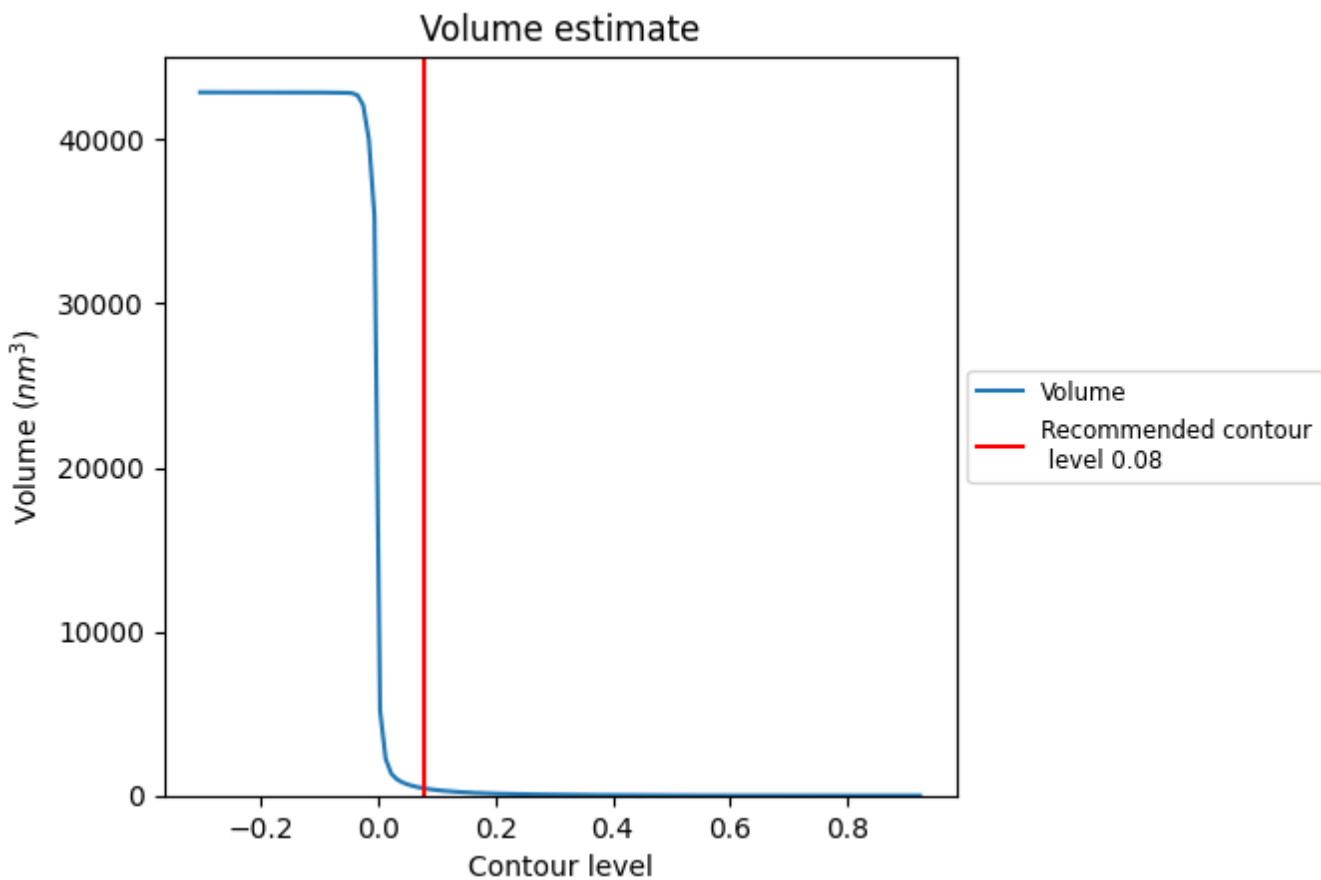
This section contains the results of statistical analysis of the map.

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.

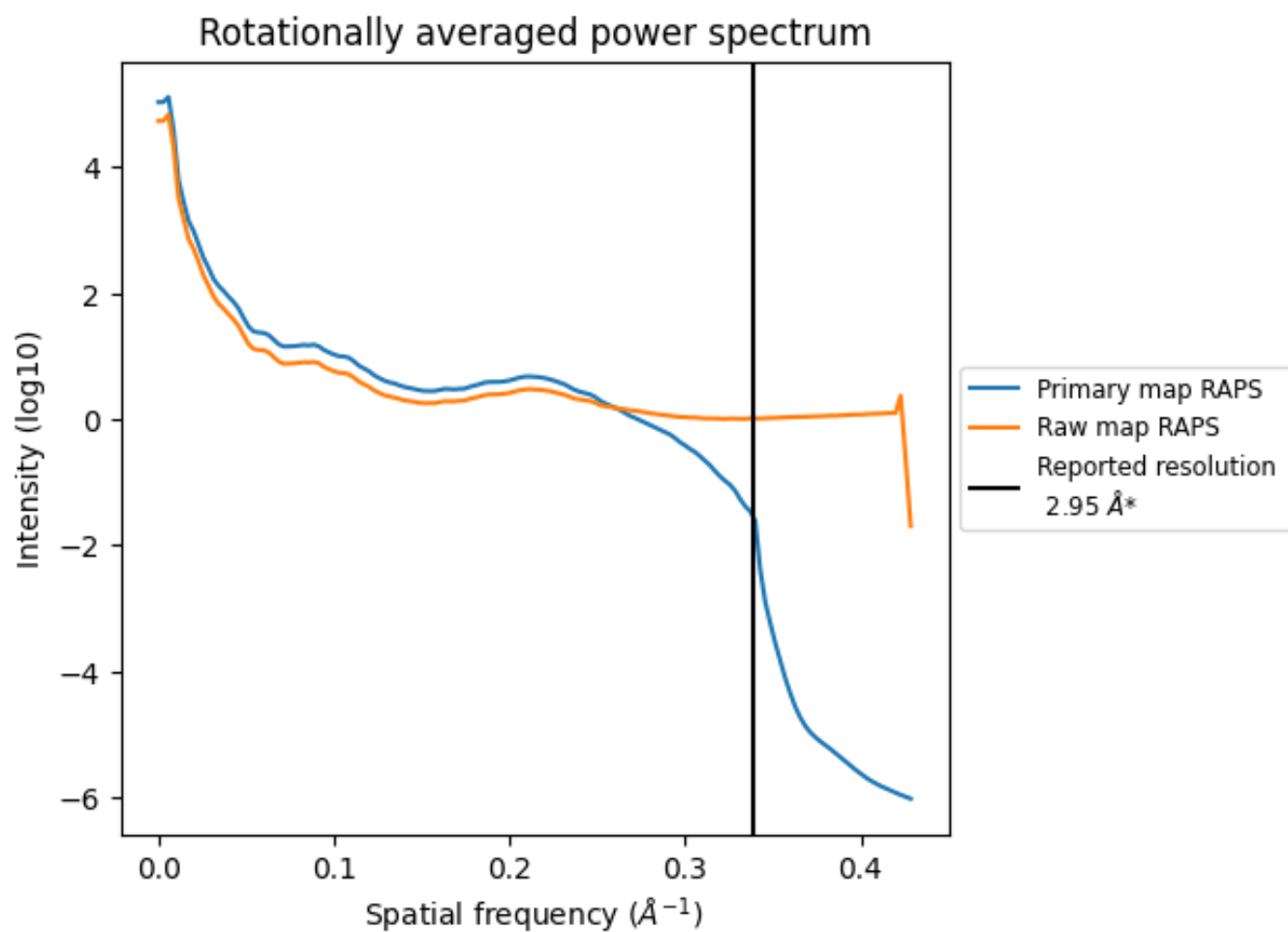
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 432 nm^3 ; this corresponds to an approximate mass of 390 kDa.

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

7.3 Rotationally averaged power spectrum i

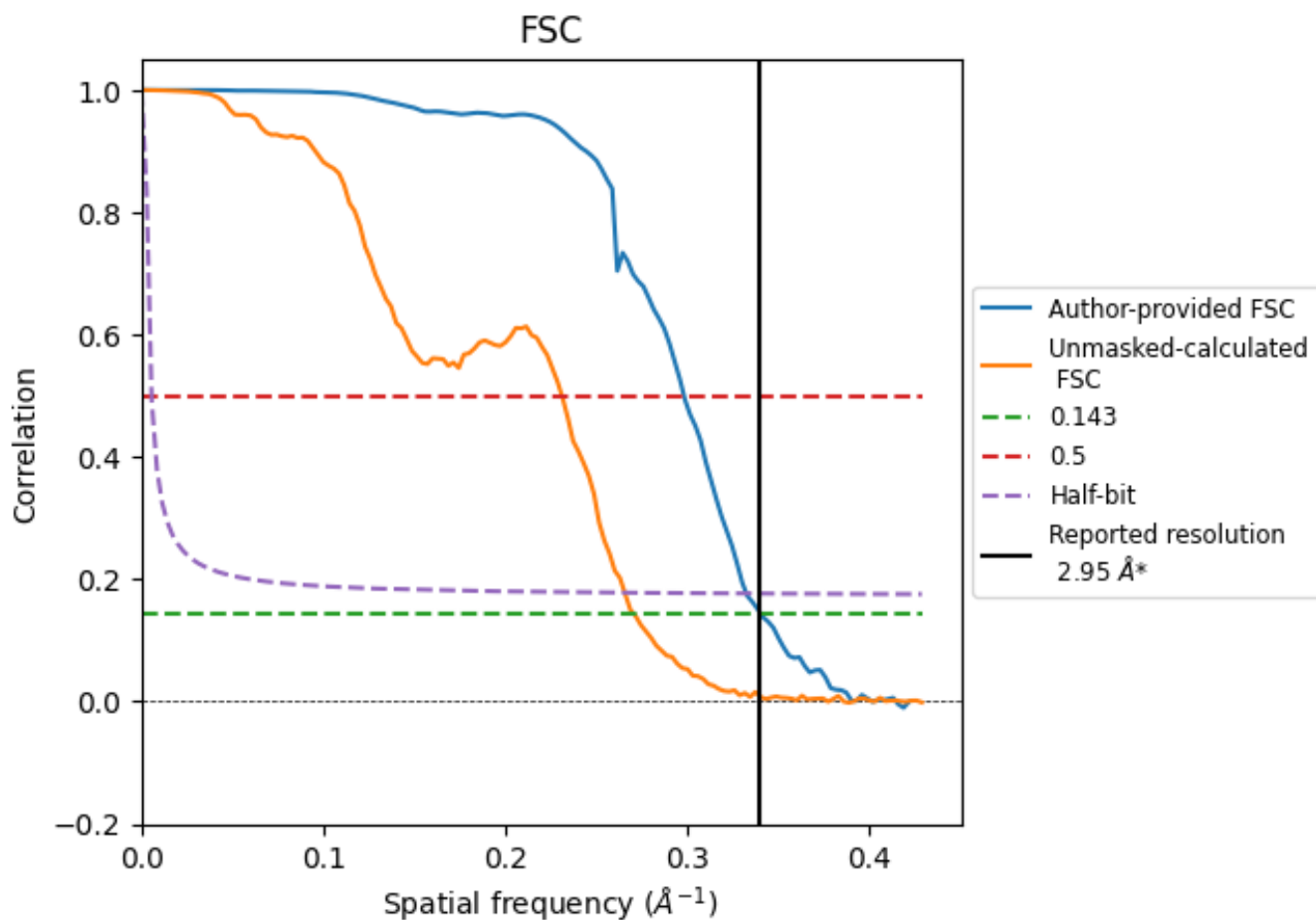


*Reported resolution corresponds to spatial frequency of 0.339 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.339 \AA^{-1}

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.95	-	-
Author-provided FSC curve	2.94	3.36	3.01
Unmasked-calculated*	3.70	4.33	3.77

*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 3.70 differs from the reported value 2.95 by more than 10 %

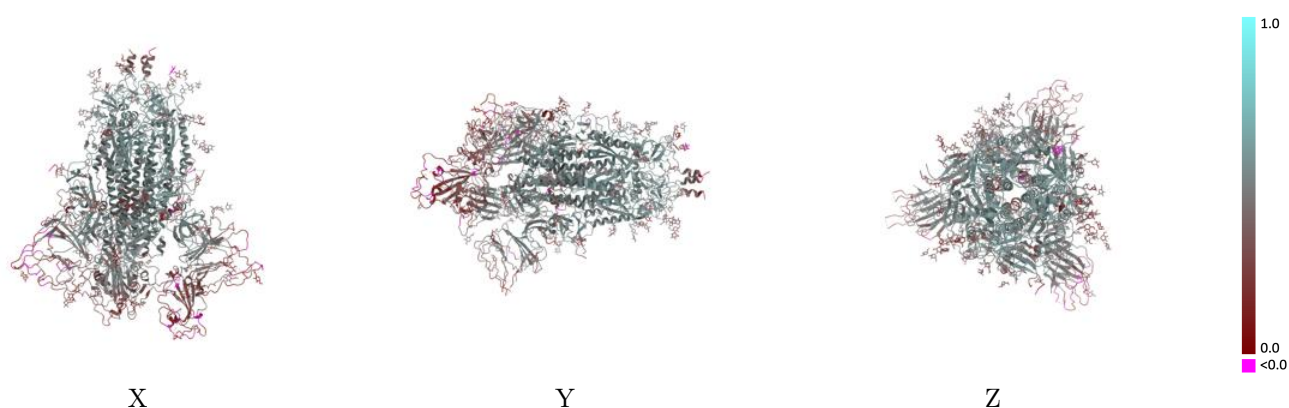
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-44447 and PDB model 9BD9. Per-residue inclusion information can be found in section 3 on page 14.

9.1 Map-model overlay [i](#)

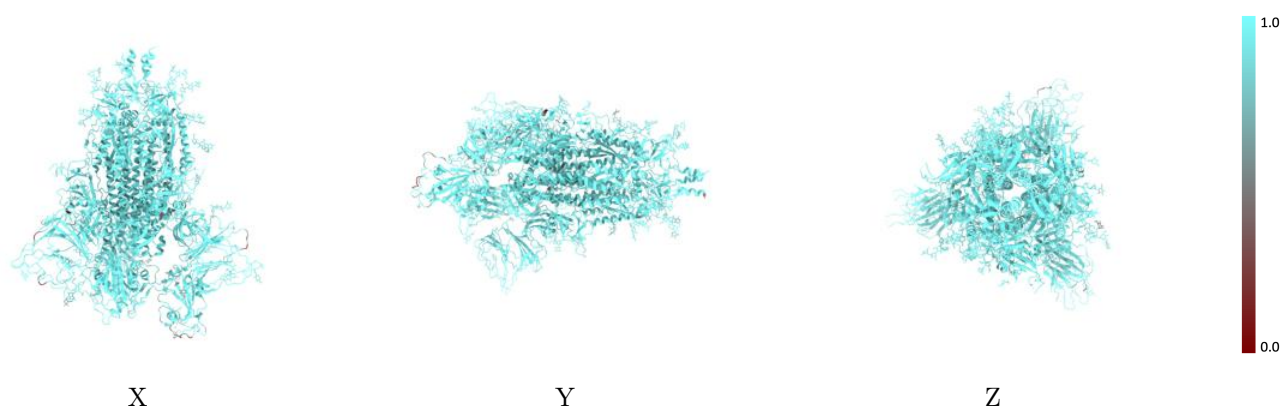
This section was not generated.

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



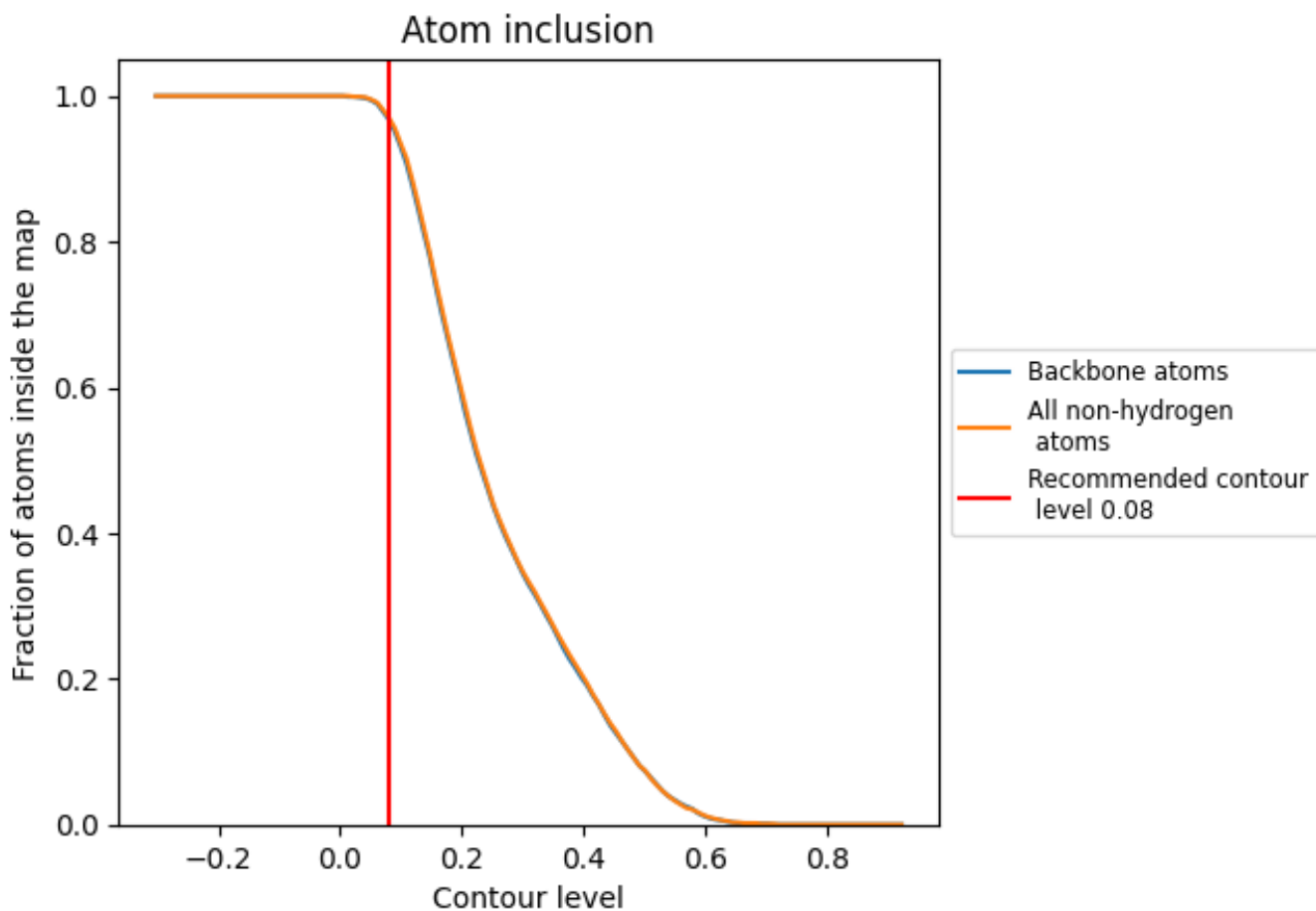
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.08).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 97% of all backbone atoms, 97% 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 (0.08) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9710	 0.4470
A	 0.9730	 0.4480
B	 0.9800	 0.4580
C	 0.9750	 0.4550
D	 0.9200	 0.3150
E	 0.8210	 0.2270
F	 0.8570	 0.2860
G	 0.8970	 0.1270
H	 0.9230	 0.2680
I	 0.9640	 0.2960
J	 1.0000	 0.4800
K	 0.8800	 0.3450
L	 0.8950	 0.2660
M	 0.9020	 0.3530
N	 0.9740	 0.4280
O	 0.8210	 0.2380
P	 0.9230	 0.2570
Q	 0.9290	 0.3840
R	 0.7860	 0.2590
S	 0.9640	 0.2880
T	 0.9230	 0.3670
U	 0.9230	 0.2870
V	 0.9600	 0.3690
W	 0.9200	 0.3710
X	 0.9210	 0.3950
Y	 0.9340	 0.2700
Z	 0.9010	 0.3180
a	 0.8210	 0.2890
b	 0.9290	 0.3910
c	 0.7140	 0.2500
d	 0.7500	 0.2590
e	 0.9400	 0.3150
f	 0.9640	 0.2680
g	 1.0000	 0.5000
h	 0.8200	 0.2610



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Chain	Atom inclusion	Q-score
i	 0.9740	 0.4340
j	 0.9400	 0.3780
k	 0.8330	 0.2780
l	 0.8000	 0.1860