



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 05:15 AM UTC

PDB ID : 6JX7 / pdb_00006jx7
EMDB ID : EMD-9891
Title : Cryo-EM structure of spike protein of feline infectious peritonitis virus strain UU4
Authors : Hsu, S.T.D.; Yang, T.J.; Ko, T.P.; Draczkowski, P.
Deposited on : 2019-04-22
Resolution : 3.31 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

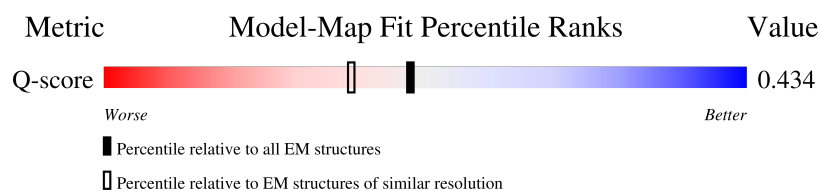
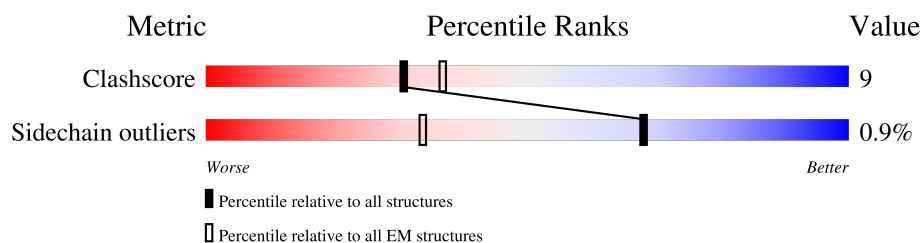
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.31 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	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14550 (2.81 - 3.81)

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	1468	16% 68% 17% 15%
1	B	1468	16% 68% 17% 15%
1	C	1468	16% 68% 17% 15%
2	0	2	50% 50% 50%
2	4	2	50% 100%

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Mol	Chain	Length	Quality of chain
2	D	2	50% 100%
2	F	2	50% 50%
2	K	2	100% 100%
2	M	2	50% 100%
2	N	2	50% 50%
2	Q	2	50% 50%
2	U	2	50% 100%
2	V	2	50% 100%
2	X	2	50% 50%
2	c	2	100% 100%
2	e	2	50% 100%
2	f	2	100% 50%
2	i	2	50% 100%
2	m	2	50% 100%
2	n	2	50% 100%
2	p	2	50% 50%
2	u	2	100% 100%
2	w	2	50% 50%
2	x	2	50% 50%
3	E	3	67% 100%
3	G	3	33% 67%
3	H	3	100% 100%
3	W	3	67% 100%
3	Y	3	33% 67%
3	Z	3	100% 100%

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Mol	Chain	Length	Quality of chain
3	o	3	67% 100%
3	q	3	67% 33% 67%
3	r	3	100% 100%
4	I	7	71% 29% 43% 29%
4	a	7	71% 29% 43% 29%
4	s	7	71% 29% 43% 29%
5	J	4	75% 100%
5	L	4	75% 75% 25%
5	O	4	75% 75% 25%
5	P	4	75% 75% 25%
5	b	4	75% 100%
5	d	4	75% 75% 25%
5	g	4	75% 50% 50%
5	h	4	75% 75% 25%
5	t	4	75% 100%
5	v	4	75% 75% 25%
5	y	4	75% 50% 50%
5	z	4	75% 75% 25%
6	1	4	75% 25% 50% 25%
6	R	4	75% 25% 50% 25%
6	j	4	75% 25% 50% 25%
7	2	5	60% 60% 40%
7	S	5	60% 60% 40%
7	k	5	60% 60% 40%
8	3	7	71% 71% 29%

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Mol	Chain	Length	Quality of chain	
8	T	7	71%	
			71% 29%	
8	1	7	71%	
			71% 29%	

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 32115 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Feline Infectious Peritonitis Virus Spike Protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1245	Total	C	N	O	S	0	0
			9768	6227	1601	1888	52		
1	B	1245	Total	C	N	O	S	0	0
			9768	6227	1601	1888	52		
1	C	1245	Total	C	N	O	S	0	0
			9768	6227	1601	1888	52		

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



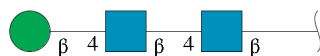
Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	2	Total	C	N	O	0	0
			28	16	2	10		
2	F	2	Total	C	N	O	0	0
			28	16	2	10		
2	K	2	Total	C	N	O	0	0
			28	16	2	10		
2	M	2	Total	C	N	O	0	0
			28	16	2	10		
2	N	2	Total	C	N	O	0	0
			28	16	2	10		
2	Q	2	Total	C	N	O	0	0
			28	16	2	10		
2	U	2	Total	C	N	O	0	0
			28	16	2	10		
2	V	2	Total	C	N	O	0	0
			28	16	2	10		
2	X	2	Total	C	N	O	0	0
			28	16	2	10		

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Mol	Chain	Residues	Atoms				AltConf	Trace
2	c	2	Total	C	N	O	0	0
			28	16	2	10		
2	e	2	Total	C	N	O	0	0
			28	16	2	10		
2	f	2	Total	C	N	O	0	0
			28	16	2	10		
2	i	2	Total	C	N	O	0	0
			28	16	2	10		
2	m	2	Total	C	N	O	0	0
			28	16	2	10		
2	n	2	Total	C	N	O	0	0
			28	16	2	10		
2	p	2	Total	C	N	O	0	0
			28	16	2	10		
2	u	2	Total	C	N	O	0	0
			28	16	2	10		
2	w	2	Total	C	N	O	0	0
			28	16	2	10		
2	x	2	Total	C	N	O	0	0
			28	16	2	10		
2	0	2	Total	C	N	O	0	0
			28	16	2	10		
2	4	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 3 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.



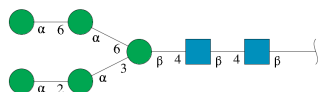
Mol	Chain	Residues	Atoms				AltConf	Trace
3	E	3	Total	C	N	O	0	0
			39	22	2	15		
3	G	3	Total	C	N	O	0	0
			39	22	2	15		
3	H	3	Total	C	N	O	0	0
			39	22	2	15		
3	W	3	Total	C	N	O	0	0
			39	22	2	15		
3	Y	3	Total	C	N	O	0	0
			39	22	2	15		

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Mol	Chain	Residues	Atoms				AltConf	Trace
3	Z	3	Total	C	N	O	0	0
			39	22	2	15		
3	o	3	Total	C	N	O	0	0
			39	22	2	15		
3	q	3	Total	C	N	O	0	0
			39	22	2	15		
3	r	3	Total	C	N	O	0	0
			39	22	2	15		

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



Mol	Chain	Residues	Atoms				AltConf	Trace
4	I	7	Total	C	N	O	0	0
			83	46	2	35		
4	a	7	Total	C	N	O	0	0
			83	46	2	35		
4	s	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 5 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.



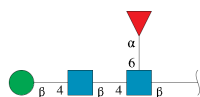
Mol	Chain	Residues	Atoms				AltConf	Trace
5	J	4	Total	C	N	O	0	0
			50	28	2	20		
5	L	4	Total	C	N	O	0	0
			50	28	2	20		
5	O	4	Total	C	N	O	0	0
			50	28	2	20		

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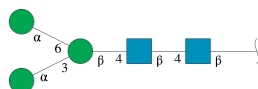
Mol	Chain	Residues	Atoms				AltConf	Trace
5	P	4	Total	C	N	O	0	0
			50	28	2	20		
5	b	4	Total	C	N	O	0	0
			50	28	2	20		
5	d	4	Total	C	N	O	0	0
			50	28	2	20		
5	g	4	Total	C	N	O	0	0
			50	28	2	20		
5	h	4	Total	C	N	O	0	0
			50	28	2	20		
5	t	4	Total	C	N	O	0	0
			50	28	2	20		
5	v	4	Total	C	N	O	0	0
			50	28	2	20		
5	y	4	Total	C	N	O	0	0
			50	28	2	20		
5	z	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 6 is an oligosaccharide called 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.



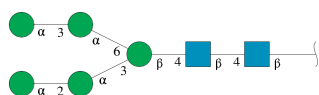
Mol	Chain	Residues	Atoms				AltConf	Trace
6	R	4	Total	C	N	O	0	0
			49	28	2	19		
6	j	4	Total	C	N	O	0	0
			49	28	2	19		
6	1	4	Total	C	N	O	0	0
			49	28	2	19		

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



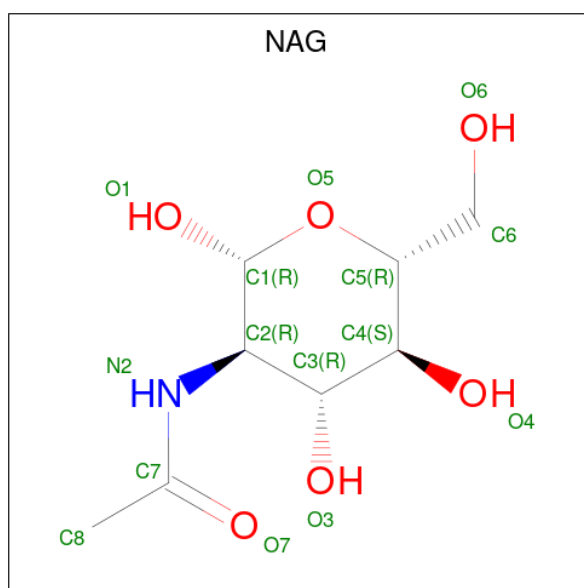
Mol	Chain	Residues	Atoms				AltConf	Trace
7	S	5	Total	C	N	O	0	0
			61	34	2	25		
7	k	5	Total	C	N	O	0	0
			61	34	2	25		
7	2	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 8 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[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	T	7	Total	C	N	O	0	0
			83	46	2	35		
8	1	7	Total	C	N	O	0	0
			83	46	2	35		
8	3	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



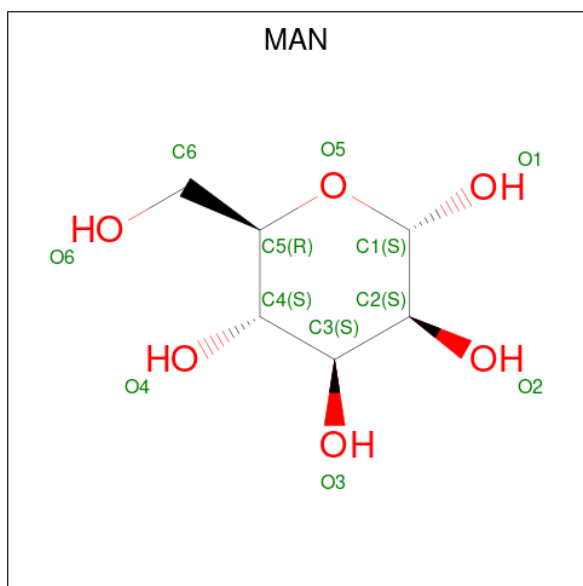
Mol	Chain	Residues	Atoms				AltConf
9	A	1	Total	C	N	O	0
			14	8	1	5	
9	A	1	Total	C	N	O	0
			14	8	1	5	
9	A	1	Total	C	N	O	0
			14	8	1	5	
9	A	1	Total	C	N	O	0
			14	8	1	5	
9	A	1	Total	C	N	O	0
			14	8	1	5	
9	A	1	Total	C	N	O	0
			14	8	1	5	
9	A	1	Total	C	N	O	0
			14	8	1	5	
9	A	1	Total	C	N	O	0
			14	8	1	5	
9	B	1	Total	C	N	O	0
			14	8	1	5	
9	B	1	Total	C	N	O	0
			14	8	1	5	
9	B	1	Total	C	N	O	0
			14	8	1	5	
9	B	1	Total	C	N	O	0
			14	8	1	5	
9	B	1	Total	C	N	O	0
			14	8	1	5	
9	B	1	Total	C	N	O	0
			14	8	1	5	
9	B	1	Total	C	N	O	0
			14	8	1	5	
9	B	1	Total	C	N	O	0
			14	8	1	5	
9	C	1	Total	C	N	O	0
			14	8	1	5	
9	C	1	Total	C	N	O	0
			14	8	1	5	
9	C	1	Total	C	N	O	0
			14	8	1	5	
9	C	1	Total	C	N	O	0
			14	8	1	5	

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

- Molecule 10 is alpha-D-mannopyranose (CCD ID: MAN) (formula: C₆H₁₂O₆).

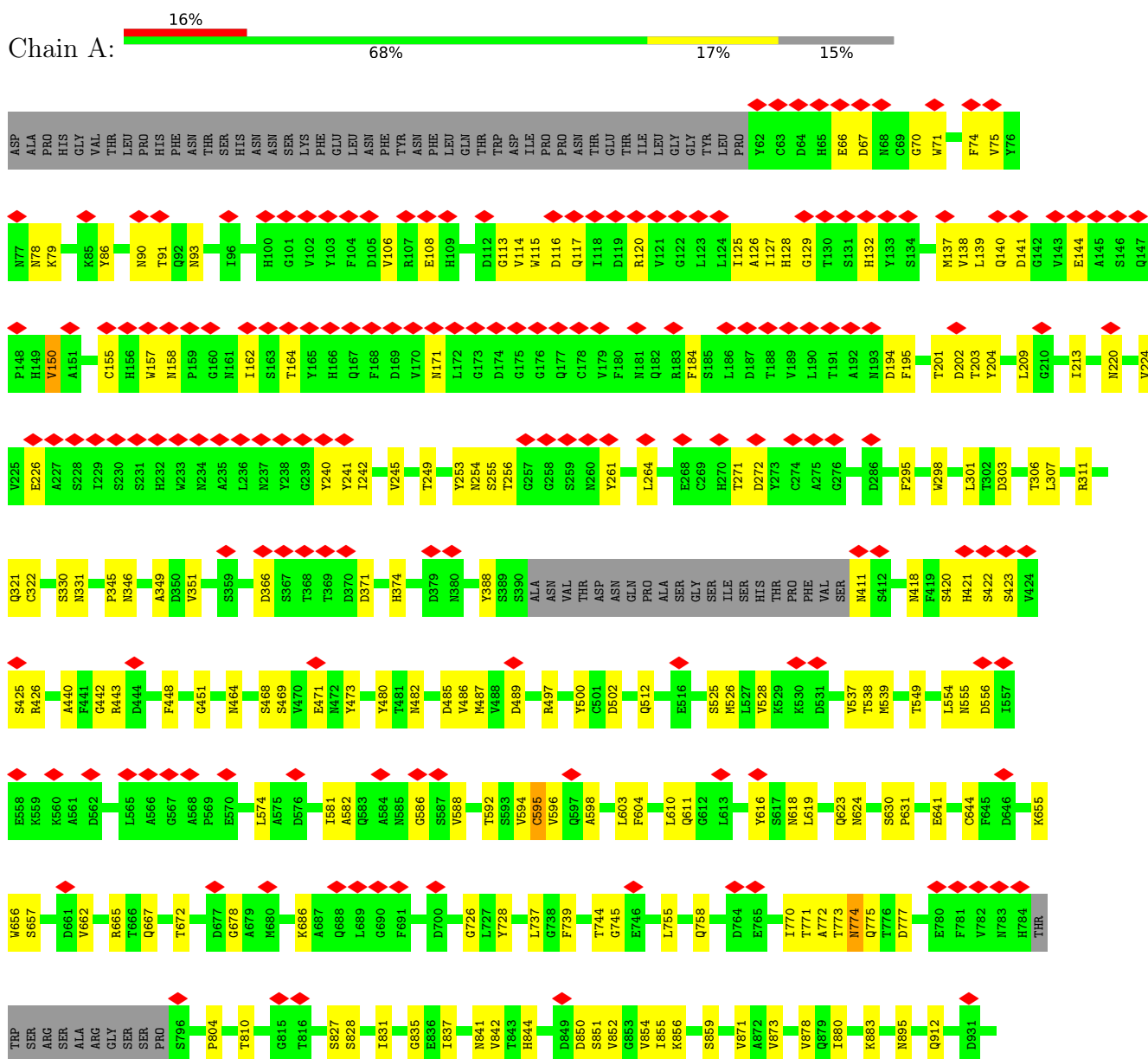


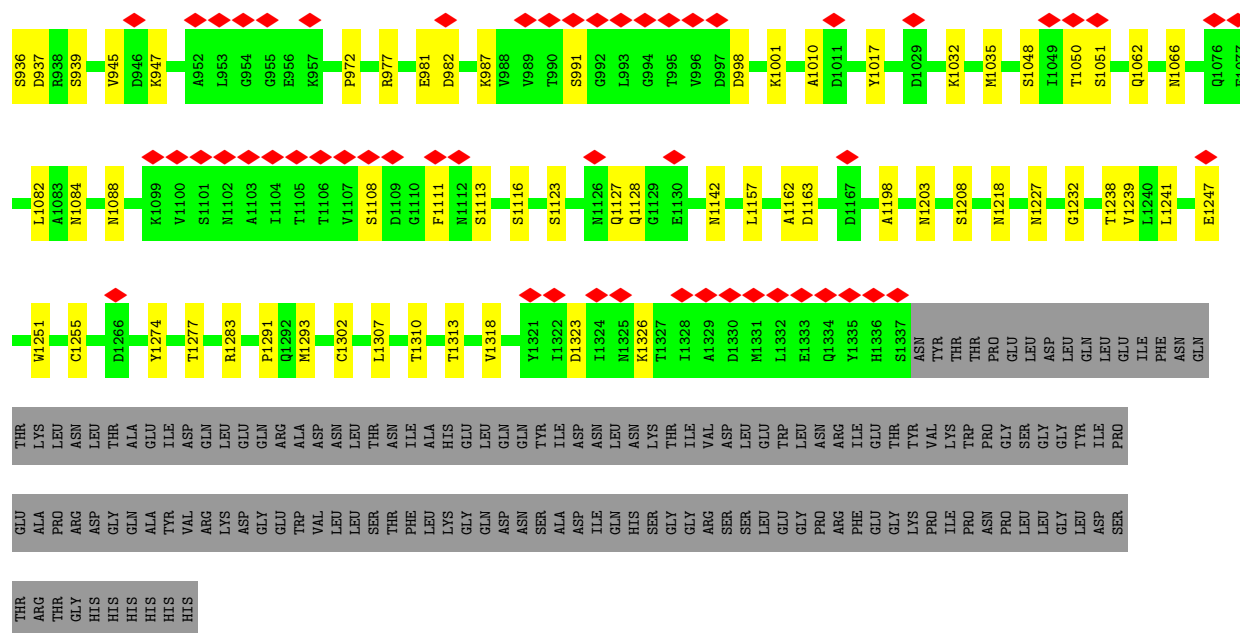
Mol	Chain	Residues	Atoms			AltConf
10	A	1	Total	C	O	0
			11	6	5	
10	A	1	Total	C	O	0
			11	6	5	
10	B	1	Total	C	O	0
			11	6	5	
10	B	1	Total	C	O	0
			11	6	5	
10	C	1	Total	C	O	0
			11	6	5	
10	C	1	Total	C	O	0
			11	6	5	

3 Residue-property plots

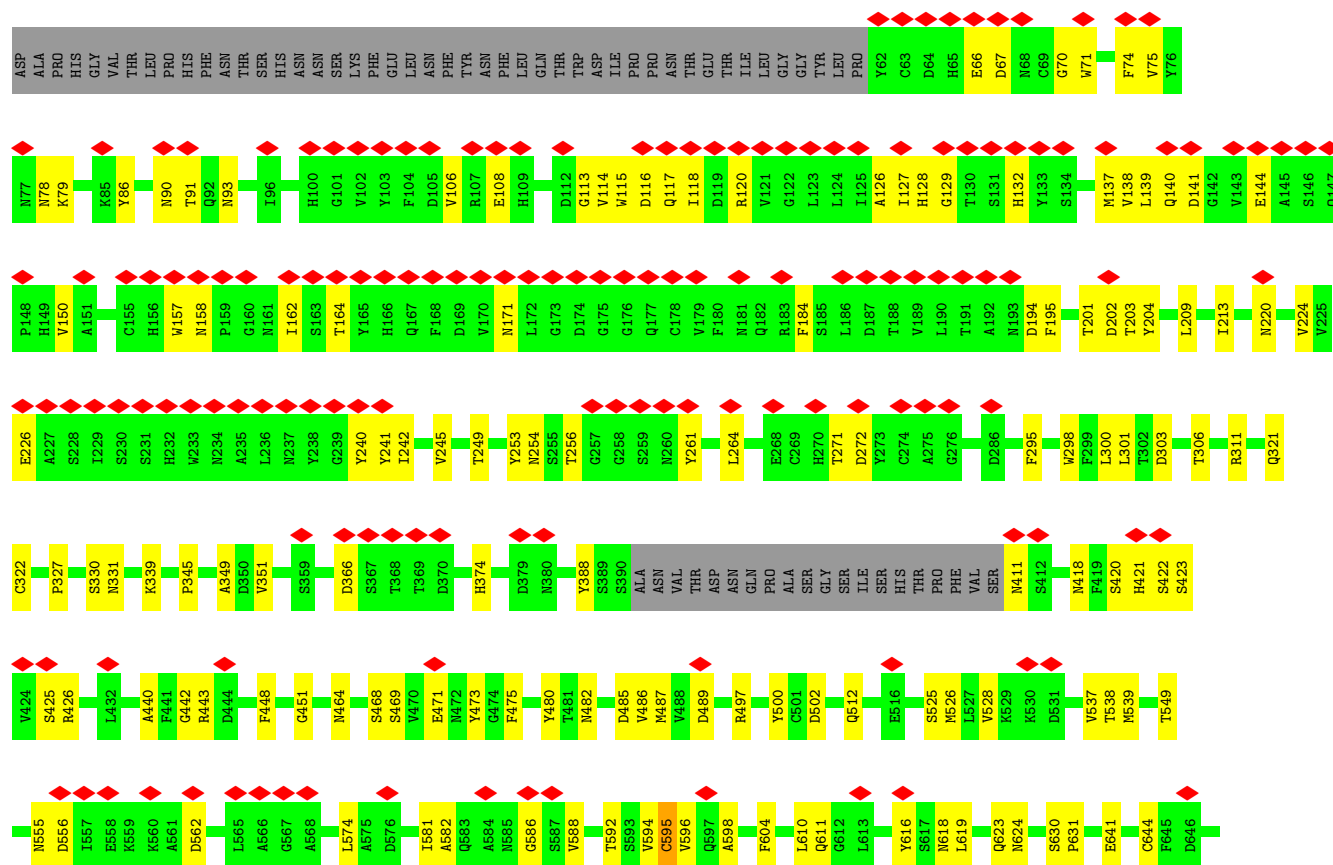
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: Feline Infectious Peritonitis Virus Spike Protein





• Molecule 1: Feline Infectious Peritonitis Virus Spike Protein



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



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



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



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



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



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



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



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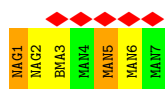
- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



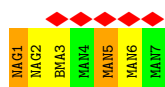
- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



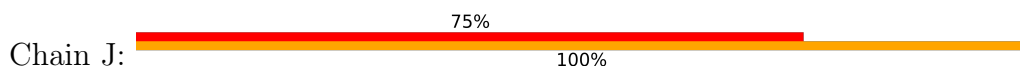
- Molecule 4: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



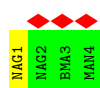
- Molecule 4: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



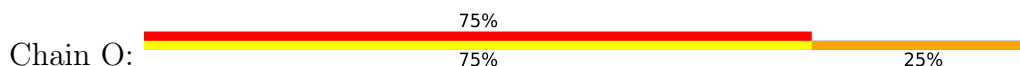
- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

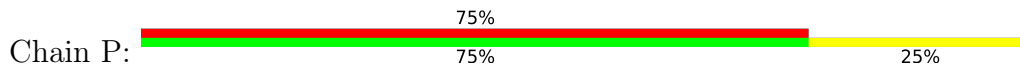


- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

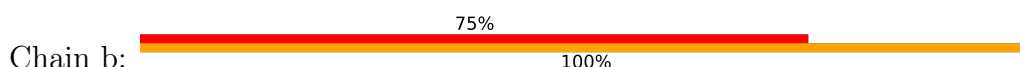




- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



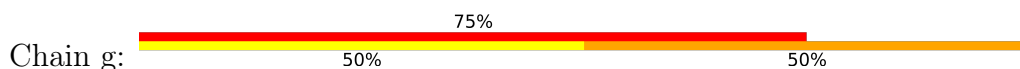
- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



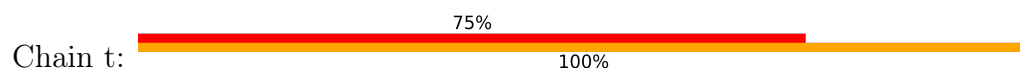
- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



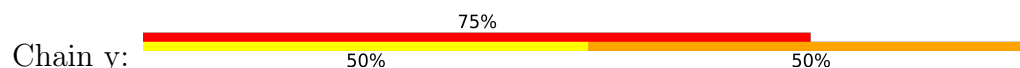
- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



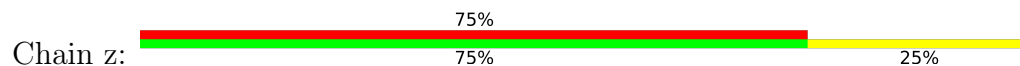
- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



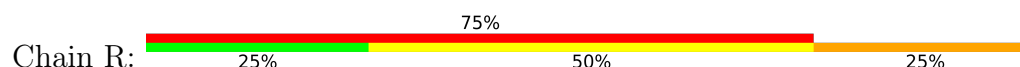
- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



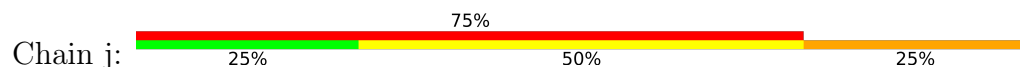
- Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



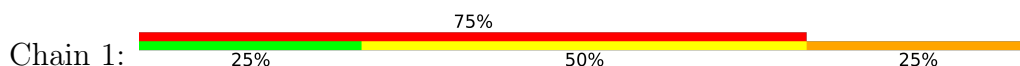
- Molecule 6: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



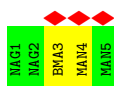
- Molecule 6: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



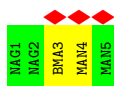
- Molecule 6: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



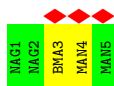
- Molecule 7: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



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



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



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



- Molecule 8: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-3)- α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	102586	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1.5	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	31.221	Depositor
Minimum map value	-16.757	Depositor
Average map value	-0.008	Depositor
Map value standard deviation	0.895	Depositor
Recommended contour level	5	Depositor
Map size (Å)	382.8, 382.8, 382.8	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.86999995, 0.86999995, 0.86999995	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.42	0/9993	0.61	3/13627 (0.0%)
1	B	0.42	0/9993	0.61	3/13627 (0.0%)
1	C	0.42	0/9993	0.61	3/13627 (0.0%)
All	All	0.42	0/29979	0.61	9/40881 (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	3
1	B	0	3
1	C	0	3
All	All	0	9

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	775	GLN	CA-C-N	-7.41	110.35	120.87
1	C	775	GLN	C-N-CA	-7.41	110.35	120.87
1	A	775	GLN	CA-C-N	-7.41	110.35	120.87
1	A	775	GLN	C-N-CA	-7.41	110.35	120.87
1	B	775	GLN	CA-C-N	-7.39	110.37	120.87
1	B	775	GLN	C-N-CA	-7.39	110.37	120.87
1	A	774	ASN	CB-CA-C	-7.38	95.72	110.42
1	B	774	ASN	CB-CA-C	-7.38	95.73	110.42
1	C	774	ASN	CB-CA-C	-7.38	95.74	110.42

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	171	ASN	Peptide
1	A	203	THR	Peptide
1	A	261	TYR	Peptide
1	B	171	ASN	Peptide
1	B	203	THR	Peptide
1	B	261	TYR	Peptide
1	C	171	ASN	Peptide
1	C	203	THR	Peptide
1	C	261	TYR	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	9768	0	9391	170	0
1	B	9768	0	9391	175	0
1	C	9768	0	9391	169	0
2	0	28	0	25	0	0
2	4	28	0	25	0	0
2	D	28	0	25	3	0
2	F	28	0	25	4	0
2	K	28	0	25	0	0
2	M	28	0	25	3	0
2	N	28	0	25	0	0
2	Q	28	0	25	0	0
2	U	28	0	25	0	0
2	V	28	0	25	2	0
2	X	28	0	25	3	0
2	c	28	0	25	0	0
2	e	28	0	25	2	0
2	f	28	0	25	0	0
2	i	28	0	25	0	0
2	m	28	0	25	0	0
2	n	28	0	25	2	0
2	p	28	0	25	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	u	28	0	25	0	0
2	w	28	0	25	2	0
2	x	28	0	25	0	0
3	E	39	0	34	0	0
3	G	39	0	34	7	0
3	H	39	0	34	6	0
3	W	39	0	34	0	0
3	Y	39	0	34	7	0
3	Z	39	0	34	6	0
3	o	39	0	34	0	0
3	q	39	0	34	7	0
3	r	39	0	34	5	0
4	I	83	0	70	3	0
4	a	83	0	70	3	0
4	s	83	0	70	3	0
5	J	50	0	43	3	0
5	L	50	0	43	0	0
5	O	50	0	43	2	0
5	P	50	0	43	0	0
5	b	50	0	43	3	0
5	d	50	0	43	0	0
5	g	50	0	43	4	0
5	h	50	0	43	0	0
5	t	50	0	43	4	0
5	v	50	0	43	0	0
5	y	50	0	43	3	0
5	z	50	0	43	0	0
6	l	49	0	43	4	0
6	R	49	0	43	4	0
6	j	49	0	43	3	0
7	2	61	0	52	0	0
7	S	61	0	52	0	0
7	k	61	0	52	0	0
8	3	83	0	70	4	0
8	T	83	0	70	4	0
8	l	83	0	70	4	0
9	A	126	0	117	2	0
9	B	126	0	117	2	0
9	C	126	0	117	2	0
10	A	22	0	20	1	0
10	B	22	0	20	1	0
10	C	22	0	20	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	32115	0	30636	560	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 9.

All (560) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:345:PRO:HB3	3:H:1:NAG:O7	1.22	1.29
1:B:345:PRO:HB3	3:Z:1:NAG:O7	1.22	1.28
1:C:345:PRO:HB3	3:r:1:NAG:O7	1.22	1.27
1:C:345:PRO:CB	3:r:1:NAG:O7	1.95	1.15
1:B:345:PRO:CB	3:Z:1:NAG:O7	1.95	1.15
1:A:345:PRO:CB	3:H:1:NAG:O7	1.95	1.13
6:1:2:NAG:H82	6:1:4:FUC:C1	1.94	0.98
6:R:2:NAG:H82	6:R:4:FUC:C1	1.94	0.97
6:j:2:NAG:H82	6:j:4:FUC:C1	1.94	0.97
3:G:2:NAG:H3	3:G:2:NAG:C8	2.17	0.75
3:Y:2:NAG:H3	3:Y:2:NAG:C8	2.17	0.75
3:q:2:NAG:C8	3:q:2:NAG:H3	2.17	0.74
1:A:854:VAL:HG12	1:A:855:ILE:N	2.04	0.73
1:C:854:VAL:HG12	1:C:855:ILE:N	2.04	0.71
1:B:854:VAL:HG12	1:B:855:ILE:N	2.04	0.70
2:M:1:NAG:O7	2:M:1:NAG:O3	2.09	0.70
2:e:1:NAG:O7	2:e:1:NAG:O3	2.09	0.70
3:G:2:NAG:H3	3:G:2:NAG:H82	1.74	0.70
3:Y:2:NAG:H3	3:Y:2:NAG:H82	1.74	0.70
3:q:2:NAG:H3	3:q:2:NAG:H82	1.74	0.69
1:A:883:LYS:HD2	1:A:1128:GLN:OE1	1.94	0.68
1:B:883:LYS:HD2	1:B:1128:GLN:OE1	1.94	0.67
1:C:883:LYS:HD2	1:C:1128:GLN:OE1	1.94	0.67
5:b:3:BMA:H4	5:b:4:MAN:H5	1.77	0.67
2:w:1:NAG:O7	2:w:1:NAG:O3	2.09	0.67
10:A:1573:MAN:H3	5:g:3:BMA:H4	1.77	0.67
5:t:3:BMA:H4	5:t:4:MAN:H5	1.77	0.66
1:A:440:ALA:HB3	1:A:448:PHE:HB2	1.77	0.66
10:C:1501:MAN:H3	5:O:3:BMA:H4	1.77	0.66
1:C:440:ALA:HB3	1:C:448:PHE:HB2	1.77	0.66
5:J:3:BMA:H4	5:J:4:MAN:H5	1.77	0.66
1:B:440:ALA:HB3	1:B:448:PHE:HB2	1.77	0.65
6:1:2:NAG:C8	6:1:4:FUC:C1	2.74	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:345:PRO:CA	3:H:1:NAG:O7	2.45	0.65
10:B:1573:MAN:H3	5:y:3:BMA:H4	1.77	0.64
1:A:202:ASP:OD2	3:G:1:NAG:H81	1.98	0.64
1:A:374:HIS:HB3	1:A:464:ASN:HB3	1.79	0.64
1:B:374:HIS:HB3	1:B:464:ASN:HB3	1.79	0.64
1:B:202:ASP:OD2	3:Y:1:NAG:H81	1.98	0.64
6:R:2:NAG:C8	6:R:4:FUC:C1	2.74	0.64
1:B:345:PRO:CA	3:Z:1:NAG:O7	2.45	0.64
1:A:253:TYR:CE1	1:A:264:LEU:HD22	2.34	0.63
1:C:374:HIS:HB3	1:C:464:ASN:HB3	1.79	0.63
1:C:345:PRO:CA	3:r:1:NAG:O7	2.45	0.63
1:B:253:TYR:CE1	1:B:264:LEU:HD22	2.34	0.63
1:C:253:TYR:CE1	1:C:264:LEU:HD22	2.34	0.63
1:C:202:ASP:OD2	3:q:1:NAG:H81	1.98	0.62
1:B:90:ASN:ND2	1:B:249:THR:O	2.33	0.62
1:A:90:ASN:ND2	1:A:249:THR:O	2.33	0.62
6:j:2:NAG:C8	6:j:4:FUC:C1	2.74	0.62
1:C:90:ASN:ND2	1:C:249:THR:O	2.33	0.61
6:1:2:NAG:C8	6:1:4:FUC:O5	2.49	0.61
1:B:883:LYS:H	1:B:1128:GLN:NE2	1.99	0.61
6:R:2:NAG:C8	6:R:4:FUC:O5	2.49	0.61
1:A:883:LYS:H	1:A:1128:GLN:NE2	1.99	0.60
1:C:883:LYS:H	1:C:1128:GLN:NE2	1.99	0.60
1:A:120:ARG:HB3	1:A:144:GLU:HG2	1.83	0.60
1:A:388:TYR:OH	1:A:426:ARG:NH2	2.35	0.60
1:B:854:VAL:CG1	1:B:855:ILE:N	2.65	0.60
6:j:2:NAG:C8	6:j:4:FUC:O5	2.49	0.60
1:C:854:VAL:CG1	1:C:855:ILE:N	2.65	0.60
1:C:388:TYR:OH	1:C:426:ARG:NH2	2.35	0.60
1:B:120:ARG:HB3	1:B:144:GLU:HG2	1.83	0.60
1:A:86:TYR:CD1	2:F:1:NAG:H82	2.37	0.60
1:B:388:TYR:OH	1:B:426:ARG:NH2	2.35	0.60
1:B:204:TYR:H	1:B:220:ASN:HD21	1.50	0.60
1:C:86:TYR:CD1	2:p:1:NAG:H82	2.37	0.60
1:A:204:TYR:H	1:A:220:ASN:HD21	1.50	0.59
1:A:202:ASP:OD2	3:G:1:NAG:C8	2.51	0.59
1:B:468:SER:OG	1:B:469:SER:N	2.35	0.59
1:A:468:SER:OG	1:A:469:SER:N	2.35	0.59
1:B:86:TYR:CD1	2:X:1:NAG:H82	2.37	0.59
1:C:883:LYS:HD2	1:C:1128:GLN:CD	2.28	0.59
1:C:120:ARG:HB3	1:C:144:GLU:HG2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:TYR:H	1:C:220:ASN:HD21	1.50	0.59
1:A:486:VAL:HG12	1:A:500:TYR:HA	1.85	0.59
1:C:202:ASP:OD2	3:q:1:NAG:C8	2.51	0.59
1:A:873:VAL:HG22	1:A:1238:THR:HG22	1.85	0.59
1:B:202:ASP:OD2	3:Y:1:NAG:C8	2.51	0.59
1:A:854:VAL:CG1	1:A:855:ILE:N	2.65	0.58
1:C:486:VAL:HG12	1:C:500:TYR:HA	1.85	0.58
1:C:873:VAL:HG22	1:C:1238:THR:HG22	1.85	0.58
1:B:883:LYS:HD2	1:B:1128:GLN:CD	2.28	0.58
1:B:486:VAL:HG12	1:B:500:TYR:HA	1.85	0.58
1:B:556:ASP:O	1:B:611:GLN:NE2	2.37	0.58
1:B:873:VAL:HG22	1:B:1238:THR:HG22	1.85	0.58
1:C:468:SER:OG	1:C:469:SER:N	2.35	0.58
1:A:556:ASP:O	1:A:611:GLN:NE2	2.37	0.58
1:A:883:LYS:HD2	1:A:1128:GLN:CD	2.28	0.58
1:A:726:GLY:H	1:C:306:THR:HG21	1.69	0.58
1:A:592:THR:HG23	9:A:1536:NAG:H81	1.86	0.57
1:A:1293:MET:SD	1:A:1293:MET:N	2.77	0.57
8:l:1:NAG:O7	8:l:1:NAG:H3	2.03	0.57
1:B:306:THR:HG21	1:C:726:GLY:H	1.68	0.57
1:B:1293:MET:SD	1:B:1293:MET:N	2.77	0.57
1:A:306:THR:HG21	1:B:726:GLY:H	1.69	0.57
1:C:592:THR:HG23	9:C:1537:NAG:H81	1.86	0.57
8:T:1:NAG:O7	8:T:1:NAG:H3	2.03	0.57
8:3:1:NAG:O7	8:3:1:NAG:H3	2.03	0.57
1:C:556:ASP:O	1:C:611:GLN:NE2	2.37	0.57
1:A:937:ASP:OD2	1:B:844:HIS:ND1	2.38	0.56
1:B:937:ASP:OD2	1:C:844:HIS:ND1	2.38	0.56
1:C:586:GLY:HA3	9:C:1536:NAG:H82	1.86	0.56
1:A:657:SER:OG	1:A:667:GLN:NE2	2.38	0.56
1:A:844:HIS:ND1	1:C:937:ASP:OD2	2.38	0.56
2:M:2:NAG:O6	2:M:2:NAG:O4	2.22	0.56
1:B:592:THR:HG23	9:B:1536:NAG:H81	1.86	0.56
1:B:945:VAL:HG12	1:B:972:PRO:HA	1.87	0.56
1:A:883:LYS:HD2	1:A:1128:GLN:NE2	2.21	0.56
1:A:586:GLY:HA3	9:A:1535:NAG:H82	1.86	0.56
1:C:1293:MET:SD	1:C:1293:MET:N	2.77	0.56
1:B:883:LYS:HD2	1:B:1128:GLN:NE2	2.21	0.56
1:C:883:LYS:HD2	1:C:1128:GLN:NE2	2.21	0.56
1:B:201:THR:OG1	1:B:202:ASP:N	2.38	0.56
1:A:1218:ASN:H	1:A:1239:VAL:HG11	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:657:SER:OG	1:C:667:GLN:NE2	2.38	0.56
1:B:586:GLY:HA3	9:B:1535:NAG:H82	1.86	0.55
1:B:1218:ASN:H	1:B:1239:VAL:HG11	1.72	0.55
1:C:1163:ASP:N	1:C:1163:ASP:OD1	2.38	0.55
4:s:5:MAN:O6	4:s:5:MAN:O4	2.19	0.55
1:A:945:VAL:HG12	1:A:972:PRO:HA	1.87	0.55
1:A:1062:GLN:O	1:A:1066:ASN:ND2	2.39	0.55
1:C:945:VAL:HG12	1:C:972:PRO:HA	1.87	0.55
1:C:1218:ASN:H	1:C:1239:VAL:HG11	1.71	0.55
1:A:91:THR:OG1	1:A:93:ASN:ND2	2.40	0.55
1:B:1062:GLN:O	1:B:1066:ASN:ND2	2.39	0.55
1:B:91:THR:OG1	1:B:93:ASN:ND2	2.40	0.55
1:B:977:ARG:NH1	1:B:982:ASP:OD1	2.40	0.55
1:C:91:THR:OG1	1:C:93:ASN:ND2	2.40	0.55
1:B:657:SER:OG	1:B:667:GLN:NE2	2.38	0.55
1:C:471:GLU:HG2	4:s:1:NAG:H81	1.89	0.55
1:C:641:GLU:HB2	1:C:678:GLY:HA3	1.89	0.54
1:A:641:GLU:HB2	1:A:678:GLY:HA3	1.89	0.54
1:B:471:GLU:HG2	4:a:1:NAG:H81	1.89	0.54
4:a:5:MAN:O6	4:a:5:MAN:O4	2.19	0.54
1:C:998:ASP:OD1	1:C:998:ASP:N	2.38	0.54
1:C:977:ARG:NH1	1:C:982:ASP:OD1	2.40	0.54
1:C:1062:GLN:O	1:C:1066:ASN:ND2	2.39	0.54
1:A:303:ASP:N	1:A:303:ASP:OD1	2.40	0.54
1:A:977:ARG:NH1	1:A:982:ASP:OD1	2.40	0.54
1:A:1241:LEU:HD21	8:T:1:NAG:O6	2.08	0.54
1:A:525:SER:OG	1:A:526:MET:N	2.41	0.54
1:A:1241:LEU:CD2	8:T:1:NAG:O6	2.56	0.54
1:B:641:GLU:HB2	1:B:678:GLY:HA3	1.89	0.54
1:C:895:ASN:OD1	1:C:895:ASN:N	2.41	0.54
5:J:1:NAG:H62	5:J:2:NAG:HN2	1.73	0.54
5:b:1:NAG:H62	5:b:2:NAG:HN2	1.73	0.54
1:A:471:GLU:HG2	4:I:1:NAG:H81	1.89	0.54
1:A:895:ASN:OD1	1:A:895:ASN:N	2.40	0.54
1:C:1241:LEU:HD21	8:3:1:NAG:O6	2.08	0.54
1:B:485:ASP:OD1	1:B:485:ASP:N	2.41	0.54
1:B:1241:LEU:CD2	8:l:1:NAG:O6	2.56	0.54
1:B:525:SER:OG	1:B:526:MET:N	2.41	0.53
1:B:1241:LEU:HD21	8:l:1:NAG:O6	2.08	0.53
1:C:303:ASP:OD1	1:C:303:ASP:N	2.40	0.53
1:A:810:THR:HG21	1:A:835:GLY:HA3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:t:1:NAG:H62	5:t:2:NAG:HN2	1.73	0.53
1:A:526:MET:HE1	1:A:777:ASP:HB2	1.91	0.53
1:B:108:GLU:HG2	1:B:224:VAL:HG13	1.90	0.53
1:C:526:MET:HE1	1:C:777:ASP:HB2	1.91	0.53
1:C:1241:LEU:CD2	8:3:1:NAG:O6	2.56	0.53
1:A:831:ILE:HG12	1:A:837:ILE:HG12	1.90	0.53
1:C:108:GLU:HG2	1:C:224:VAL:HG13	1.91	0.53
2:w:1:NAG:HO3	2:w:1:NAG:C7	2.14	0.53
1:B:1163:ASP:OD1	1:B:1163:ASP:N	2.38	0.53
1:C:201:THR:OG1	1:C:202:ASP:N	2.38	0.53
1:C:525:SER:OG	1:C:526:MET:N	2.41	0.53
1:C:810:THR:HG21	1:C:835:GLY:HA3	1.91	0.53
1:C:330:SER:OG	1:C:331:ASN:N	2.41	0.53
1:B:810:THR:HG21	1:B:835:GLY:HA3	1.91	0.53
1:C:485:ASP:OD1	1:C:485:ASP:N	2.41	0.53
1:A:108:GLU:HG2	1:A:224:VAL:HG13	1.91	0.52
1:A:841:ASN:HD22	5:O:1:NAG:H83	1.74	0.52
1:B:831:ILE:HG12	1:B:837:ILE:HG12	1.90	0.52
1:B:330:SER:OG	1:B:331:ASN:N	2.41	0.52
1:C:831:ILE:HG12	1:C:837:ILE:HG12	1.90	0.52
1:B:303:ASP:OD1	1:B:303:ASP:N	2.40	0.52
1:B:526:MET:HE1	1:B:777:ASP:HB2	1.91	0.52
1:B:895:ASN:N	1:B:895:ASN:OD1	2.40	0.52
1:B:841:ASN:HD22	5:g:1:NAG:H83	1.74	0.52
1:B:880:ILE:HG12	1:B:1232:GLY:HA2	1.92	0.52
3:G:2:NAG:C8	3:G:2:NAG:C3	2.85	0.52
1:A:349:ALA:O	1:A:443:ARG:NH2	2.43	0.52
1:A:411:ASN:OD1	1:A:411:ASN:N	2.43	0.52
1:B:549:THR:HG22	1:B:604:PHE:HB2	1.92	0.51
1:B:883:LYS:HB2	1:B:1128:GLN:NE2	2.25	0.51
1:C:883:LYS:HB2	1:C:1128:GLN:NE2	2.25	0.51
1:A:981:GLU:OE2	1:A:1227:ASN:ND2	2.44	0.51
1:A:1032:LYS:HE3	1:B:842:VAL:HB	1.92	0.51
1:A:330:SER:OG	1:A:331:ASN:N	2.41	0.51
1:A:549:THR:HG22	1:A:604:PHE:HB2	1.92	0.51
1:B:349:ALA:O	1:B:443:ARG:NH2	2.43	0.51
1:B:471:GLU:HG3	1:B:473:TYR:HB2	1.92	0.51
1:B:1157:LEU:HB2	1:B:1162:ALA:HB2	1.92	0.51
1:A:471:GLU:HG3	1:A:473:TYR:HB2	1.92	0.51
1:C:981:GLU:OE2	1:C:1227:ASN:ND2	2.44	0.51
1:A:366:ASP:N	1:A:366:ASP:OD1	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:883:LYS:HB2	1:A:1128:GLN:NE2	2.25	0.51
1:B:1032:LYS:HE3	1:C:842:VAL:HB	1.92	0.51
1:C:349:ALA:O	1:C:443:ARG:NH2	2.43	0.51
1:C:1123:SER:O	1:C:1127:GLN:NE2	2.42	0.51
1:B:366:ASP:OD1	1:B:366:ASP:N	2.43	0.51
1:C:411:ASN:OD1	1:C:411:ASN:N	2.43	0.51
1:C:880:ILE:HG12	1:C:1232:GLY:HA2	1.92	0.51
1:A:1157:LEU:HB2	1:A:1162:ALA:HB2	1.92	0.51
1:C:655:LYS:HA	1:C:672:THR:HA	1.93	0.51
1:A:880:ILE:HG12	1:A:1232:GLY:HA2	1.92	0.50
1:B:987:LYS:O	1:B:991:SER:OG	2.29	0.50
1:C:841:ASN:HD22	5:y:1:NAG:H83	1.74	0.50
1:B:1010:ALA:HB3	1:C:737:LEU:HD13	1.93	0.50
4:I:5:MAN:O6	4:I:5:MAN:O4	2.19	0.50
1:B:555:ASN:HA	1:B:610:LEU:HD12	1.93	0.50
1:A:842:VAL:HB	1:C:1032:LYS:HE3	1.92	0.50
1:A:987:LYS:O	1:A:991:SER:OG	2.29	0.50
1:A:1163:ASP:OD1	1:A:1163:ASP:N	2.38	0.50
1:C:549:THR:HG22	1:C:604:PHE:HB2	1.92	0.50
1:C:1157:LEU:HB2	1:C:1162:ALA:HB2	1.92	0.50
1:A:74:PHE:O	1:A:78:ASN:ND2	2.45	0.50
1:B:981:GLU:OE2	1:B:1227:ASN:ND2	2.44	0.50
1:C:74:PHE:O	1:C:78:ASN:ND2	2.45	0.50
1:A:485:ASP:N	1:A:485:ASP:OD1	2.41	0.50
1:A:655:LYS:HA	1:A:672:THR:HA	1.93	0.50
1:A:771:THR:OG1	1:A:772:ALA:N	2.45	0.50
1:A:1247:GLU:OE1	1:C:1050:THR:OG1	2.30	0.50
1:C:471:GLU:HG3	1:C:473:TYR:HB2	1.92	0.50
1:A:555:ASN:HA	1:A:610:LEU:HD12	1.93	0.50
1:B:655:LYS:HA	1:B:672:THR:HA	1.93	0.49
1:C:987:LYS:O	1:C:991:SER:OG	2.29	0.49
3:Y:2:NAG:C8	3:Y:2:NAG:C3	2.85	0.49
1:B:1123:SER:O	1:B:1127:GLN:NE2	2.42	0.49
1:B:74:PHE:O	1:B:78:ASN:ND2	2.45	0.49
1:B:1084:ASN:O	1:B:1088:ASN:ND2	2.46	0.49
1:B:126:ALA:HB3	1:B:226:GLU:HB2	1.94	0.49
1:B:771:THR:OG1	1:B:772:ALA:N	2.45	0.49
1:A:1010:ALA:HB3	1:B:737:LEU:HD13	1.93	0.49
1:C:771:THR:OG1	1:C:772:ALA:N	2.45	0.49
1:A:141:ASP:OD1	1:A:141:ASP:N	2.46	0.49
1:A:737:LEU:HD13	1:C:1010:ALA:HB3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1050:THR:OG1	1:B:1247:GLU:OE1	2.30	0.49
1:B:411:ASN:N	1:B:411:ASN:OD1	2.43	0.49
1:B:998:ASP:OD1	1:B:998:ASP:N	2.38	0.49
1:B:1050:THR:OG1	1:C:1247:GLU:OE1	2.30	0.49
1:A:86:TYR:CE1	2:F:1:NAG:C8	2.96	0.49
1:A:1084:ASN:O	1:A:1088:ASN:ND2	2.46	0.49
1:B:106:VAL:HG21	2:V:1:NAG:O4	2.13	0.49
1:B:420:SER:OG	1:B:421:HIS:N	2.45	0.49
1:C:86:TYR:CE1	2:p:1:NAG:C8	2.96	0.49
1:A:201:THR:OG1	1:A:202:ASP:N	2.38	0.48
1:C:139:LEU:HD21	1:C:157:TRP:HE1	1.78	0.48
1:A:1123:SER:O	1:A:1127:GLN:NE2	2.42	0.48
1:C:126:ALA:HB3	1:C:226:GLU:HB2	1.94	0.48
1:C:255:SER:O	1:C:255:SER:OG	2.31	0.48
1:C:555:ASN:HA	1:C:610:LEU:HD12	1.93	0.48
3:Y:1:NAG:H62	3:Y:2:NAG:C7	2.43	0.48
1:B:86:TYR:CE1	2:X:1:NAG:C8	2.96	0.48
1:A:254:ASN:OD1	1:A:256:THR:OG1	2.17	0.48
1:B:254:ASN:OD1	1:B:256:THR:OG1	2.17	0.48
1:C:420:SER:OG	1:C:421:HIS:N	2.45	0.48
1:B:139:LEU:HD21	1:B:157:TRP:HE1	1.78	0.48
1:C:1084:ASN:O	1:C:1088:ASN:ND2	2.46	0.48
3:q:1:NAG:H62	3:q:2:NAG:C7	2.43	0.48
1:A:420:SER:OG	1:A:421:HIS:N	2.45	0.48
1:B:1113:SER:O	1:B:1116:SER:OG	2.31	0.48
2:M:1:NAG:HO3	2:M:1:NAG:C7	2.20	0.48
1:A:139:LEU:HD21	1:A:157:TRP:HE1	1.78	0.48
8:l:6:MAN:H3	8:l:7:MAN:H2	1.54	0.48
1:A:106:VAL:HG21	2:D:1:NAG:O4	2.13	0.48
1:A:126:ALA:HB3	1:A:226:GLU:HB2	1.94	0.48
1:A:555:ASN:HB2	1:A:574:LEU:HD22	1.96	0.48
1:C:79:LYS:HD2	1:C:213:ILE:HD11	1.96	0.48
1:B:141:ASP:N	1:B:141:ASP:OD1	2.46	0.47
1:C:321:GLN:NE2	1:C:482:ASN:OD1	2.47	0.47
3:G:1:NAG:H62	3:G:2:NAG:C7	2.43	0.47
1:B:158:ASN:OD1	1:B:158:ASN:N	2.44	0.47
1:B:322:CYS:HB3	1:B:480:TYR:HD2	1.80	0.47
1:A:321:GLN:NE2	1:A:482:ASN:OD1	2.47	0.47
1:A:322:CYS:HB3	1:A:480:TYR:HD2	1.80	0.47
1:A:489:ASP:HB2	1:A:497:ARG:H	1.79	0.47
1:C:106:VAL:HG21	2:n:1:NAG:O4	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:489:ASP:HB2	1:C:497:ARG:H	1.79	0.47
1:B:321:GLN:NE2	1:B:482:ASN:OD1	2.47	0.47
3:q:2:NAG:C8	3:q:2:NAG:C3	2.85	0.47
1:A:79:LYS:HD2	1:A:213:ILE:HD11	1.96	0.47
1:B:79:LYS:HD2	1:B:213:ILE:HD11	1.96	0.47
1:C:859:SER:OG	1:C:1283:ARG:NH2	2.48	0.47
3:H:1:NAG:H62	3:H:2:NAG:C7	2.45	0.47
5:t:1:NAG:H82	5:t:1:NAG:H2	1.75	0.47
1:A:883:LYS:HD2	1:A:1128:GLN:HE22	1.79	0.47
1:A:1113:SER:O	1:A:1116:SER:OG	2.31	0.47
1:B:150:VAL:O	1:B:184:PHE:N	2.43	0.47
1:B:883:LYS:HD2	1:B:1128:GLN:HE22	1.79	0.47
3:H:2:NAG:H61	3:H:3:BMA:H2	1.97	0.47
3:Z:1:NAG:H62	3:Z:2:NAG:C7	2.45	0.47
1:C:883:LYS:HD2	1:C:1128:GLN:HE22	1.79	0.47
1:A:859:SER:OG	1:A:1283:ARG:NH2	2.48	0.46
1:B:489:ASP:HB2	1:B:497:ARG:H	1.79	0.46
1:B:1017:TYR:HE2	1:C:755:LEU:HD13	1.80	0.46
1:C:295:PHE:HD2	1:C:298:TRP:HB2	1.81	0.46
1:C:322:CYS:HB3	1:C:480:TYR:HD2	1.80	0.46
3:r:2:NAG:H61	3:r:3:BMA:H2	1.97	0.46
1:A:623:GLN:HE21	1:A:665:ARG:HH11	1.62	0.46
3:r:1:NAG:H62	3:r:2:NAG:C7	2.45	0.46
1:A:1255:CYS:HB2	1:A:1302:CYS:HB2	1.43	0.46
1:B:859:SER:OG	1:B:1283:ARG:NH2	2.48	0.46
1:C:141:ASP:OD1	1:C:141:ASP:N	2.46	0.46
1:C:204:TYR:H	1:C:220:ASN:ND2	2.14	0.46
1:A:150:VAL:O	1:A:184:PHE:N	2.43	0.46
1:A:1274:TYR:N	1:A:1277:THR:O	2.48	0.46
1:A:998:ASP:OD1	1:A:998:ASP:N	2.38	0.46
1:B:555:ASN:HB2	1:B:574:LEU:HD22	1.96	0.46
1:B:1001:LYS:NZ	1:C:804:PRO:O	2.49	0.46
1:C:623:GLN:HE21	1:C:665:ARG:HH11	1.62	0.46
1:A:936:SER:OG	1:A:939:SER:OG	2.31	0.46
8:3:6:MAN:H3	8:3:7:MAN:H2	1.54	0.46
1:A:204:TYR:H	1:A:220:ASN:ND2	2.14	0.46
1:C:1113:SER:O	1:C:1116:SER:OG	2.31	0.46
1:A:755:LEU:HD13	1:C:1017:TYR:HE2	1.80	0.46
1:B:595:CYS:HB2	1:B:644:CYS:HB3	1.57	0.46
1:C:150:VAL:HG13	1:C:184:PHE:HB2	1.98	0.46
1:A:295:PHE:HD2	1:A:298:TRP:HB2	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:PHE:HD2	1:B:298:TRP:HB2	1.81	0.45
3:Z:2:NAG:H61	3:Z:3:BMA:H2	1.97	0.45
1:A:1017:TYR:HE2	1:B:755:LEU:HD13	1.80	0.45
1:B:418:ASN:ND2	1:B:425:SER:O	2.50	0.45
1:B:1274:TYR:N	1:B:1277:THR:O	2.48	0.45
1:C:418:ASN:ND2	1:C:425:SER:O	2.50	0.45
1:C:555:ASN:HB2	1:C:574:LEU:HD22	1.96	0.45
1:A:773:THR:HG22	1:A:774:ASN:HB2	1.98	0.45
1:B:1032:LYS:HA	1:B:1035:MET:HE3	1.98	0.45
1:C:986:ASN:O	1:C:990:THR:OG1	2.33	0.45
1:C:1108:SER:H	1:C:1111:PHE:HE2	1.64	0.45
1:C:1274:TYR:N	1:C:1277:THR:O	2.48	0.45
1:B:150:VAL:HG13	1:B:184:PHE:HB2	1.98	0.45
1:B:623:GLN:HE21	1:B:665:ARG:HH11	1.62	0.45
1:C:422:SER:OG	1:C:423:SER:N	2.50	0.45
1:A:804:PRO:O	1:C:1001:LYS:NZ	2.49	0.45
1:A:1001:LYS:NZ	1:B:804:PRO:O	2.49	0.45
1:A:209:LEU:HD23	1:A:209:LEU:HA	1.86	0.45
1:A:345:PRO:HB3	3:H:1:NAG:C7	2.26	0.45
1:B:422:SER:OG	1:B:423:SER:N	2.50	0.45
1:C:595:CYS:HB2	1:C:644:CYS:HB3	1.57	0.45
1:C:150:VAL:O	1:C:184:PHE:N	2.43	0.45
1:C:744:THR:OG1	1:C:745:GLY:N	2.50	0.45
1:C:936:SER:OG	1:C:939:SER:OG	2.31	0.45
1:A:75:VAL:HG11	1:A:194:ASP:HA	1.99	0.45
1:A:150:VAL:HG13	1:A:184:PHE:HB2	1.98	0.45
1:B:75:VAL:HG11	1:B:194:ASP:HA	1.99	0.45
1:B:538:THR:OG1	1:B:539:MET:N	2.50	0.45
1:C:538:THR:OG1	1:C:539:MET:N	2.50	0.45
1:C:773:THR:HG22	1:C:774:ASN:HB2	1.98	0.45
5:b:3:BMA:H4	5:b:4:MAN:C5	2.45	0.45
1:A:603:LEU:HD12	1:A:603:LEU:HA	1.84	0.45
1:A:744:THR:OG1	1:A:745:GLY:N	2.50	0.45
1:C:1032:LYS:HA	1:C:1035:MET:HE3	1.98	0.45
5:t:3:BMA:H4	5:t:4:MAN:C5	2.45	0.45
1:A:418:ASN:ND2	1:A:425:SER:O	2.50	0.45
1:B:67:ASP:O	1:B:71:TRP:N	2.48	0.45
1:B:114:VAL:HG11	1:B:195:PHE:HB3	1.99	0.45
1:B:1142:ASN:OD1	1:B:1142:ASN:N	2.49	0.45
1:B:1174:LEU:HD23	1:B:1174:LEU:HA	1.86	0.45
2:e:2:NAG:O6	2:e:2:NAG:O4	2.22	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:371:ASP:OD1	1:A:371:ASP:N	2.48	0.44
1:A:1123:SER:OG	1:A:1127:GLN:NE2	2.50	0.44
1:A:1310:THR:HG22	1:A:1313:THR:HG22	1.99	0.44
1:B:113:GLY:HA2	1:B:245:VAL:HG23	1.99	0.44
1:B:1310:THR:HG22	1:B:1313:THR:HG22	1.99	0.44
1:C:114:VAL:HG11	1:C:195:PHE:HB3	1.99	0.44
1:C:684:MET:HE3	1:C:684:MET:HB3	1.86	0.44
1:A:624:ASN:HB2	1:A:630:SER:HA	2.00	0.44
1:A:1032:LYS:HA	1:A:1035:MET:HE3	1.98	0.44
1:B:1108:SER:H	1:B:1111:PHE:HE2	1.64	0.44
1:C:66:GLU:O	1:C:70:GLY:N	2.48	0.44
1:C:631:PRO:HG3	1:C:656:TRP:CD2	2.53	0.44
1:A:140:GLN:HE21	1:A:150:VAL:HG23	1.82	0.44
1:A:631:PRO:HG3	1:A:656:TRP:CD2	2.53	0.44
1:A:854:VAL:CG1	1:A:855:ILE:H	2.30	0.44
1:B:744:THR:OG1	1:B:745:GLY:N	2.50	0.44
1:B:947:LYS:HA	1:B:947:LYS:HD3	1.85	0.44
1:B:66:GLU:O	1:B:70:GLY:N	2.48	0.44
1:B:631:PRO:HG3	1:B:656:TRP:CD2	2.53	0.44
1:C:67:ASP:O	1:C:71:TRP:N	2.48	0.44
1:C:1123:SER:OG	1:C:1127:GLN:NE2	2.50	0.44
5:g:2:NAG:O4	5:g:2:NAG:O6	2.35	0.44
3:q:2:NAG:H3	3:q:2:NAG:H83	1.97	0.44
1:C:113:GLY:HA2	1:C:245:VAL:HG23	1.99	0.44
5:J:3:BMA:H4	5:J:4:MAN:C5	2.46	0.44
1:A:422:SER:OG	1:A:423:SER:N	2.50	0.44
1:A:538:THR:OG1	1:A:539:MET:N	2.50	0.44
1:A:1108:SER:H	1:A:1111:PHE:HE2	1.64	0.44
1:B:140:GLN:HE21	1:B:150:VAL:HG23	1.82	0.44
1:B:1123:SER:OG	1:B:1127:GLN:NE2	2.50	0.44
1:B:773:THR:HG22	1:B:774:ASN:HB2	1.98	0.44
1:B:204:TYR:H	1:B:220:ASN:ND2	2.14	0.44
1:B:502:ASP:OD1	1:B:502:ASP:N	2.50	0.44
1:B:997:ASP:OD1	1:B:997:ASP:N	2.42	0.44
1:C:1251:TRP:CE2	1:C:1307:LEU:HD13	2.53	0.44
1:A:502:ASP:N	1:A:502:ASP:OD1	2.50	0.43
1:A:1251:TRP:CE2	1:A:1307:LEU:HD13	2.53	0.43
1:A:113:GLY:HA2	1:A:245:VAL:HG23	1.99	0.43
1:A:162:ILE:HA	1:A:164:THR:HG23	2.00	0.43
1:C:75:VAL:HG11	1:C:194:ASP:HA	1.99	0.43
1:A:912:GLN:NE2	1:B:512:GLN:OE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:LEU:HD23	1:B:209:LEU:HA	1.86	0.43
1:B:912:GLN:NE2	1:C:512:GLN:OE1	2.52	0.43
1:B:1255:CYS:HB2	1:B:1302:CYS:HB2	1.43	0.43
1:A:114:VAL:HG11	1:A:195:PHE:HB3	1.99	0.43
1:A:1142:ASN:N	1:A:1142:ASN:OD1	2.49	0.43
1:B:1251:TRP:CE2	1:B:1307:LEU:HD13	2.53	0.43
1:C:502:ASP:OD1	1:C:502:ASP:N	2.50	0.43
3:Y:2:NAG:H3	3:Y:2:NAG:H83	1.97	0.43
1:A:67:ASP:O	1:A:71:TRP:N	2.48	0.43
1:B:728:TYR:HA	1:B:739:PHE:HA	2.01	0.43
1:B:986:ASN:O	1:B:990:THR:OG1	2.33	0.43
1:C:1310:THR:HG22	1:C:1313:THR:HG22	1.99	0.43
1:C:854:VAL:CG1	1:C:855:ILE:H	2.30	0.43
1:C:1142:ASN:OD1	1:C:1142:ASN:N	2.49	0.43
1:B:116:ASP:N	1:B:241:TYR:O	2.52	0.43
1:A:1203:ASN:O	1:A:1208:SER:OG	2.33	0.43
1:B:581:ILE:HD12	1:B:594:VAL:HG21	2.01	0.43
1:B:582:ALA:HA	1:B:588:VAL:HG23	2.01	0.43
1:C:603:LEU:HD12	1:C:603:LEU:HA	1.84	0.43
4:s:1:NAG:H82	4:s:1:NAG:H2	1.76	0.43
1:A:595:CYS:HB2	1:A:644:CYS:HB3	1.57	0.43
1:A:883:LYS:CD	1:A:1128:GLN:HE22	2.32	0.43
1:B:562:ASP:OD1	1:B:562:ASP:N	2.52	0.43
1:C:140:GLN:HE21	1:C:150:VAL:HG23	1.82	0.43
1:C:624:ASN:HB2	1:C:630:SER:HA	2.00	0.43
8:T:6:MAN:H3	8:T:7:MAN:H2	1.54	0.43
1:B:856:LYS:HD3	1:B:856:LYS:HA	1.91	0.42
1:A:311:ARG:HH11	1:A:487:MET:HE1	1.84	0.42
1:A:554:LEU:HD23	1:A:554:LEU:HA	1.89	0.42
1:A:610:LEU:HD23	1:A:610:LEU:HA	1.88	0.42
1:A:758:GLN:HB2	1:A:770:ILE:HB	2.01	0.42
1:B:758:GLN:HB2	1:B:770:ILE:HB	2.01	0.42
1:C:162:ILE:HA	1:C:164:THR:HG23	2.00	0.42
1:C:311:ARG:HH11	1:C:487:MET:HE1	1.84	0.42
1:A:526:MET:HE3	1:A:526:MET:HB2	1.92	0.42
1:B:311:ARG:HH11	1:B:487:MET:HE1	1.84	0.42
1:B:624:ASN:HB2	1:B:630:SER:HA	2.00	0.42
1:C:301:LEU:HD23	1:C:451:GLY:HA2	2.01	0.42
1:A:351:VAL:HG12	1:A:442:GLY:HA2	2.01	0.42
1:C:1291:PRO:HG2	1:C:1318:VAL:HA	2.01	0.42
1:A:947:LYS:HD3	1:A:947:LYS:HA	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:854:VAL:CG1	1:B:855:ILE:H	2.30	0.42
1:B:984:LEU:HD23	1:B:984:LEU:HA	1.90	0.42
1:C:116:ASP:N	1:C:241:TYR:O	2.52	0.42
1:C:158:ASN:OD1	1:C:158:ASN:N	2.44	0.42
1:C:758:GLN:HB2	1:C:770:ILE:HB	2.01	0.42
1:C:1326:LYS:HE2	1:C:1326:LYS:HB3	1.91	0.42
1:A:115:TRP:HA	1:A:242:ILE:HA	2.01	0.42
1:A:116:ASP:N	1:A:241:TYR:O	2.52	0.42
1:A:827:SER:OG	1:A:828:SER:N	2.52	0.42
1:B:127:ILE:HB	1:B:138:VAL:HB	2.02	0.42
1:C:618:ASN:N	1:C:618:ASN:OD1	2.52	0.42
1:C:728:TYR:HA	1:C:739:PHE:HA	2.01	0.42
1:C:1174:LEU:HD23	1:C:1174:LEU:HA	1.86	0.42
1:A:850:ASP:O	1:A:852:VAL:N	2.48	0.42
1:B:301:LEU:HD23	1:B:451:GLY:HA2	2.01	0.42
1:C:127:ILE:HB	1:C:138:VAL:HB	2.02	0.42
1:C:351:VAL:HG12	1:C:442:GLY:HA2	2.01	0.42
1:C:883:LYS:CD	1:C:1128:GLN:HE22	2.32	0.42
1:C:1241:LEU:HD23	1:C:1241:LEU:HA	1.88	0.42
1:B:827:SER:OG	1:B:828:SER:N	2.52	0.42
1:B:883:LYS:HA	1:B:884:PRO:HD3	1.90	0.42
1:B:1291:PRO:HG2	1:B:1318:VAL:HA	2.01	0.42
1:C:353:ARG:NH2	1:C:438:GLU:OE2	2.41	0.42
1:C:616:TYR:CE1	1:C:662:VAL:HG13	2.55	0.42
1:C:927:LEU:HD23	1:C:927:LEU:HA	1.86	0.42
1:A:128:HIS:ND1	1:A:129:GLY:O	2.53	0.42
1:A:255:SER:O	1:A:255:SER:OG	2.31	0.42
1:B:128:HIS:ND1	1:B:129:GLY:O	2.53	0.42
1:B:850:ASP:O	1:B:852:VAL:N	2.48	0.42
1:C:1332:LEU:HD23	1:C:1332:LEU:HA	1.86	0.42
6:R:2:NAG:H82	6:R:4:FUC:O5	2.15	0.42
1:B:883:LYS:H	1:B:1128:GLN:HE21	1.67	0.42
2:V:2:NAG:O7	2:V:2:NAG:H3	2.20	0.42
1:A:871:VAL:HG13	1:A:1082:LEU:HD22	2.02	0.41
1:B:115:TRP:HA	1:B:242:ILE:HA	2.01	0.41
1:B:162:ILE:HA	1:B:164:THR:HG23	2.00	0.41
1:C:115:TRP:HA	1:C:242:ILE:HA	2.01	0.41
1:C:582:ALA:HA	1:C:588:VAL:HG23	2.01	0.41
4:I:1:NAG:H82	4:I:1:NAG:H2	1.76	0.41
1:A:512:GLN:OE1	1:C:912:GLN:NE2	2.52	0.41
1:A:582:ALA:HA	1:A:588:VAL:HG23	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:616:TYR:CE1	1:A:662:VAL:HG13	2.55	0.41
1:A:686:LYS:HE3	1:A:686:LYS:HB2	1.86	0.41
1:A:728:TYR:HA	1:A:739:PHE:HA	2.01	0.41
1:A:773:THR:HG22	1:A:774:ASN:N	2.36	0.41
1:C:581:ILE:HD12	1:C:594:VAL:HG21	2.01	0.41
1:C:1323:ASP:HB3	1:C:1326:LYS:HE3	2.02	0.41
1:A:581:ILE:HD12	1:A:594:VAL:HG21	2.01	0.41
1:B:117:GLN:HB2	1:B:240:TYR:HD1	1.86	0.41
1:B:339:LYS:HE2	1:B:339:LYS:HB3	1.87	0.41
1:B:883:LYS:CD	1:B:1128:GLN:HE22	2.32	0.41
1:C:366:ASP:OD1	1:C:366:ASP:N	2.43	0.41
2:D:2:NAG:O7	2:D:2:NAG:H3	2.20	0.41
1:A:66:GLU:O	1:A:70:GLY:N	2.48	0.41
1:A:1291:PRO:HG2	1:A:1318:VAL:HA	2.01	0.41
1:C:86:TYR:CZ	2:p:1:NAG:C8	3.04	0.41
2:n:2:NAG:O7	2:n:2:NAG:H3	2.20	0.41
1:A:117:GLN:HB2	1:A:240:TYR:HD1	1.86	0.41
1:A:618:ASN:N	1:A:618:ASN:OD1	2.52	0.41
1:A:1048:SER:OG	1:A:1051:SER:OG	2.36	0.41
1:B:351:VAL:HG12	1:B:442:GLY:HA2	2.01	0.41
1:B:616:TYR:CE1	1:B:662:VAL:HG13	2.55	0.41
1:B:618:ASN:OD1	1:B:618:ASN:N	2.52	0.41
1:C:117:GLN:HB2	1:C:240:TYR:HD1	1.86	0.41
1:C:1111:PHE:O	1:C:1115:ALA:N	2.45	0.41
1:C:1136:ILE:HD13	1:C:1136:ILE:HA	1.92	0.41
4:a:1:NAG:H82	4:a:1:NAG:H2	1.76	0.41
1:A:127:ILE:HB	1:A:138:VAL:HB	2.02	0.41
1:A:301:LEU:HD23	1:A:451:GLY:HA2	2.01	0.41
1:B:1323:ASP:HB3	1:B:1326:LYS:HE3	2.03	0.41
1:C:773:THR:HG22	1:C:774:ASN:N	2.36	0.41
1:C:827:SER:OG	1:C:828:SER:N	2.52	0.41
1:C:128:HIS:ND1	1:C:129:GLY:O	2.53	0.41
1:C:390:SER:O	1:C:390:SER:OG	2.33	0.41
1:C:984:LEU:HD23	1:C:984:LEU:HA	1.90	0.41
1:A:307:LEU:HD23	1:A:307:LEU:HA	1.91	0.41
1:A:596:VAL:HG22	1:A:598:ALA:H	1.86	0.41
1:B:86:TYR:CZ	2:X:1:NAG:C8	3.04	0.41
1:B:271:THR:OG1	1:B:272:ASP:N	2.54	0.41
1:B:773:THR:HG22	1:B:774:ASN:N	2.36	0.41
1:B:1222:LEU:HD23	1:B:1222:LEU:HA	1.88	0.41
1:C:209:LEU:HD23	1:C:209:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:271:THR:OG1	1:C:272:ASP:N	2.54	0.41
1:C:314:SER:HG	1:C:316:GLN:HE21	1.62	0.41
5:g:2:NAG:H4	5:g:3:BMA:H2	1.72	0.41
1:C:1163:ASP:HA	1:C:1166:VAL:HG12	2.03	0.41
1:A:86:TYR:CZ	2:F:1:NAG:C8	3.04	0.40
1:A:132:HIS:HD2	1:A:137:MET:HE1	1.87	0.40
1:A:271:THR:OG1	1:A:272:ASP:N	2.54	0.40
1:A:851:SER:HB3	1:A:856:LYS:NZ	2.36	0.40
1:B:596:VAL:HG22	1:B:598:ALA:H	1.86	0.40
1:B:132:HIS:HD2	1:B:137:MET:HE1	1.86	0.40
1:B:300:LEU:HD23	1:B:300:LEU:HA	1.94	0.40
1:B:619:LEU:HD23	1:B:619:LEU:HA	1.91	0.40
1:B:1111:PHE:O	1:B:1115:ALA:N	2.45	0.40
1:C:1255:CYS:HB2	1:C:1302:CYS:HB2	1.43	0.40
1:A:86:TYR:CG	2:F:1:NAG:H82	2.57	0.40
1:A:878:VAL:HG11	1:A:1198:ALA:HB2	2.04	0.40
1:A:1323:ASP:HB3	1:A:1326:LYS:HE3	2.02	0.40
1:B:117:GLN:OE1	1:B:118:ILE:N	2.54	0.40
1:B:1040:LEU:HD23	1:B:1040:LEU:HA	1.96	0.40
1:C:315:SER:O	1:C:315:SER:OG	2.36	0.40
3:G:2:NAG:H3	3:G:2:NAG:H83	1.97	0.40
1:B:327:PRO:HA	1:B:475:PHE:HD1	1.86	0.40
1:B:684:MET:HE3	1:B:684:MET:HB3	1.86	0.40
1:C:871:VAL:HG13	1:C:1082:LEU:HD22	2.02	0.40
5:y:2:NAG:H4	5:y:3:BMA:H2	1.72	0.40
1:A:106:VAL:HG21	2:D:1:NAG:H5	2.04	0.40
1:A:619:LEU:HD23	1:A:619:LEU:HA	1.91	0.40
1:B:345:PRO:HB3	3:Z:1:NAG:C7	2.26	0.40
1:B:851:SER:HB3	1:B:856:LYS:NZ	2.36	0.40
1:B:927:LEU:HA	1:B:927:LEU:HD23	1.86	0.40
1:B:1059:MET:HE3	1:B:1059:MET:HB2	1.90	0.40
1:B:1139:LEU:HD23	1:B:1139:LEU:HA	1.93	0.40
1:C:686:LYS:HE3	1:C:686:LYS:HB2	1.86	0.40
1:C:851:SER:HB3	1:C:856:LYS:NZ	2.36	0.40
6:1:2:NAG:H82	6:1:4:FUC:C2	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

There are no protein backbone outliers to report in this entry.

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	469/1287 (36%)	461 (98%)	8 (2%)	53	70
1	B	548/1287 (43%)	544 (99%)	4 (1%)	76	79
1	C	622/1287 (48%)	620 (100%)	2 (0%)	86	85
All	All	1639/3861 (42%)	1625 (99%)	14 (1%)	68	77

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125	ILE
1	A	150	VAL
1	A	155	CYS
1	A	158	ASN
1	A	346	ASN
1	A	528	VAL
1	A	537	VAL
1	A	595	CYS
1	B	528	VAL
1	B	537	VAL
1	B	595	CYS
1	B	974	VAL
1	C	125	ILE
1	C	974	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	ASN
1	A	77	ASN
1	A	93	ASN
1	A	132	HIS
1	A	156	HIS
1	A	177	GLN
1	A	220	ASN
1	A	555	ASN
1	A	623	GLN
1	B	418	ASN
1	B	555	ASN
1	B	611	GLN
1	B	623	GLN
1	B	667	GLN
1	B	688	GLN
1	B	1066	ASN
1	B	1088	ASN
1	B	1122	GLN
1	B	1126	ASN
1	B	1127	GLN
1	B	1128	GLN
1	B	1144	GLN
1	B	1165	GLN
1	C	623	GLN
1	C	667	GLN
1	C	688	GLN
1	C	1066	ASN
1	C	1088	ASN
1	C	1122	GLN
1	C	1126	ASN
1	C	1127	GLN
1	C	1128	GLN
1	C	1144	GLN
1	C	1165	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

186 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	0	1	2,1	14,14,15	0.37	0	17,19,21	0.52	0
2	NAG	0	2	2	14,14,15	0.32	0	17,19,21	0.59	1 (5%)
6	NAG	1	1	6,1	14,14,15	0.32	0	17,19,21	0.88	0
6	NAG	1	2	6	14,14,15	0.65	0	17,19,21	1.89	2 (11%)
6	BMA	1	3	6	11,11,12	0.59	0	15,15,17	1.08	1 (6%)
6	FUC	1	4	6	10,10,11	0.79	0	14,14,16	0.73	0
7	NAG	2	1	7,1	14,14,15	0.32	0	17,19,21	0.59	0
7	NAG	2	2	7	14,14,15	0.21	0	17,19,21	0.57	0
7	BMA	2	3	7	11,11,12	0.58	0	15,15,17	2.36	2 (13%)
7	MAN	2	4	7	11,11,12	0.56	0	15,15,17	3.21	4 (26%)
7	MAN	2	5	7	11,11,12	0.25	0	15,15,17	0.51	0
8	NAG	3	1	8,1	14,14,15	2.09	7 (50%)	17,19,21	2.37	9 (52%)
8	NAG	3	2	8	14,14,15	0.28	0	17,19,21	0.94	1 (5%)
8	BMA	3	3	8	11,11,12	0.21	0	15,15,17	1.23	2 (13%)
8	MAN	3	4	8	11,11,12	0.26	0	15,15,17	0.94	1 (6%)
8	MAN	3	5	8	11,11,12	0.28	0	15,15,17	1.18	3 (20%)
8	MAN	3	6	8	11,11,12	0.31	0	15,15,17	0.84	0
8	MAN	3	7	8	11,11,12	0.97	1 (9%)	15,15,17	0.96	2 (13%)
2	NAG	4	1	2,1	14,14,15	0.18	0	17,19,21	0.59	0
2	NAG	4	2	2	14,14,15	0.27	0	17,19,21	0.52	0
2	NAG	D	1	2,1	14,14,15	0.29	0	17,19,21	0.57	0
2	NAG	D	2	2	14,14,15	0.29	0	17,19,21	0.82	0
3	NAG	E	1	3,1	14,14,15	0.45	0	17,19,21	2.00	6 (35%)
3	NAG	E	2	3	14,14,15	0.48	0	17,19,21	1.39	2 (11%)
3	BMA	E	3	3	11,11,12	0.44	0	15,15,17	1.36	2 (13%)
2	NAG	F	1	2,1	14,14,15	0.32	0	17,19,21	0.67	0
2	NAG	F	2	2	14,14,15	0.32	0	17,19,21	0.89	0
3	NAG	G	1	3,1	14,14,15	0.33	0	17,19,21	0.54	0
3	NAG	G	2	3	14,14,15	0.25	0	17,19,21	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BMA	G	3	3	11,11,12	0.23	0	15,15,17	0.76	0
3	NAG	H	1	3,1	14,14,15	0.29	0	17,19,21	0.85	0
3	NAG	H	2	3	14,14,15	0.27	0	17,19,21	0.58	0
3	BMA	H	3	3	11,11,12	0.24	0	15,15,17	0.75	0
4	NAG	I	1	4,1	14,14,15	0.68	0	17,19,21	1.94	4 (23%)
4	NAG	I	2	4	14,14,15	0.50	0	17,19,21	1.68	4 (23%)
4	BMA	I	3	4	11,11,12	0.61	0	15,15,17	2.23	8 (53%)
4	MAN	I	4	4	11,11,12	0.69	0	15,15,17	0.96	0
4	MAN	I	5	4	11,11,12	0.60	0	15,15,17	1.38	2 (13%)
4	MAN	I	6	4	11,11,12	0.66	0	15,15,17	1.23	1 (6%)
4	MAN	I	7	4	11,11,12	0.67	0	15,15,17	0.94	0
5	NAG	J	1	5,1	14,14,15	0.54	0	17,19,21	1.27	3 (17%)
5	NAG	J	2	5	14,14,15	1.06	1 (7%)	17,19,21	1.85	3 (17%)
5	BMA	J	3	5	11,11,12	0.59	0	15,15,17	1.42	2 (13%)
5	MAN	J	4	5	11,11,12	0.38	0	15,15,17	1.23	2 (13%)
2	NAG	K	1	2,1	14,14,15	0.25	0	17,19,21	0.63	0
2	NAG	K	2	2	14,14,15	0.21	0	17,19,21	0.62	0
5	NAG	L	1	5,1	14,14,15	0.36	0	17,19,21	0.76	1 (5%)
5	NAG	L	2	5	14,14,15	0.27	0	17,19,21	0.65	0
5	BMA	L	3	5	11,11,12	0.24	0	15,15,17	0.73	0
5	MAN	L	4	5	11,11,12	0.24	0	15,15,17	0.73	0
2	NAG	M	1	2,1	14,14,15	1.25	2 (14%)	17,19,21	2.63	7 (41%)
2	NAG	M	2	2	14,14,15	0.51	0	17,19,21	1.16	2 (11%)
2	NAG	N	1	2,1	14,14,15	0.18	0	17,19,21	0.72	1 (5%)
2	NAG	N	2	2	14,14,15	0.34	0	17,19,21	0.54	0
5	NAG	O	1	5,1	14,14,15	0.27	0	17,19,21	0.65	0
5	NAG	O	2	5	14,14,15	1.04	1 (7%)	17,19,21	1.97	4 (23%)
5	BMA	O	3	5	11,11,12	0.35	0	15,15,17	1.90	4 (26%)
5	MAN	O	4	5	11,11,12	0.41	0	15,15,17	0.92	1 (6%)
5	NAG	P	1	5,1	14,14,15	0.18	0	17,19,21	0.74	1 (5%)
5	NAG	P	2	5	14,14,15	0.24	0	17,19,21	0.57	0
5	BMA	P	3	5	11,11,12	0.25	0	15,15,17	0.52	0
5	MAN	P	4	5	11,11,12	0.28	0	15,15,17	0.52	0
2	NAG	Q	1	2,1	14,14,15	0.38	0	17,19,21	0.52	0
2	NAG	Q	2	2	14,14,15	0.32	0	17,19,21	0.59	1 (5%)
6	NAG	R	1	6,1	14,14,15	0.32	0	17,19,21	0.88	0
6	NAG	R	2	6	14,14,15	0.65	0	17,19,21	1.89	2 (11%)
6	BMA	R	3	6	11,11,12	0.59	0	15,15,17	1.08	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	FUC	R	4	6	10,10,11	0.79	0	14,14,16	0.73	0
7	NAG	S	1	7,1	14,14,15	0.32	0	17,19,21	0.59	0
7	NAG	S	2	7	14,14,15	0.22	0	17,19,21	0.57	0
7	BMA	S	3	7	11,11,12	0.58	0	15,15,17	2.36	2 (13%)
7	MAN	S	4	7	11,11,12	0.56	0	15,15,17	3.21	4 (26%)
7	MAN	S	5	7	11,11,12	0.26	0	15,15,17	0.51	0
8	NAG	T	1	8,1	14,14,15	2.10	7 (50%)	17,19,21	2.37	9 (52%)
8	NAG	T	2	8	14,14,15	0.29	0	17,19,21	0.95	1 (5%)
8	BMA	T	3	8	11,11,12	0.22	0	15,15,17	1.23	2 (13%)
8	MAN	T	4	8	11,11,12	0.26	0	15,15,17	0.95	1 (6%)
8	MAN	T	5	8	11,11,12	0.29	0	15,15,17	1.18	3 (20%)
8	MAN	T	6	8	11,11,12	0.31	0	15,15,17	0.85	0
8	MAN	T	7	8	11,11,12	0.97	1 (9%)	15,15,17	0.96	2 (13%)
2	NAG	U	1	2,1	14,14,15	0.17	0	17,19,21	0.59	0
2	NAG	U	2	2	14,14,15	0.27	0	17,19,21	0.52	0
2	NAG	V	1	2,1	14,14,15	0.29	0	17,19,21	0.56	0
2	NAG	V	2	2	14,14,15	0.29	0	17,19,21	0.81	0
3	NAG	W	1	3,1	14,14,15	0.45	0	17,19,21	2.00	6 (35%)
3	NAG	W	2	3	14,14,15	0.49	0	17,19,21	1.38	2 (11%)
3	BMA	W	3	3	11,11,12	0.43	0	15,15,17	1.36	2 (13%)
2	NAG	X	1	2,1	14,14,15	0.32	0	17,19,21	0.67	0
2	NAG	X	2	2	14,14,15	0.32	0	17,19,21	0.89	0
3	NAG	Y	1	3,1	14,14,15	0.33	0	17,19,21	0.54	0
3	NAG	Y	2	3	14,14,15	0.25	0	17,19,21	0.58	0
3	BMA	Y	3	3	11,11,12	0.23	0	15,15,17	0.76	0
3	NAG	Z	1	3,1	14,14,15	0.28	0	17,19,21	0.85	0
3	NAG	Z	2	3	14,14,15	0.27	0	17,19,21	0.58	0
3	BMA	Z	3	3	11,11,12	0.23	0	15,15,17	0.75	0
4	NAG	a	1	4,1	14,14,15	0.68	0	17,19,21	1.94	4 (23%)
4	NAG	a	2	4	14,14,15	0.49	0	17,19,21	1.67	4 (23%)
4	BMA	a	3	4	11,11,12	0.60	0	15,15,17	2.24	8 (53%)
4	MAN	a	4	4	11,11,12	0.69	0	15,15,17	0.96	0
4	MAN	a	5	4	11,11,12	0.60	0	15,15,17	1.38	2 (13%)
4	MAN	a	6	4	11,11,12	0.66	0	15,15,17	1.23	1 (6%)
4	MAN	a	7	4	11,11,12	0.67	0	15,15,17	0.94	0
5	NAG	b	1	5,1	14,14,15	0.54	0	17,19,21	1.27	3 (17%)
5	NAG	b	2	5	14,14,15	1.06	1 (7%)	17,19,21	1.85	3 (17%)
5	BMA	b	3	5	11,11,12	0.59	0	15,15,17	1.42	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MAN	b	4	5	11,11,12	0.39	0	15,15,17	1.23	2 (13%)
2	NAG	c	1	2,1	14,14,15	0.24	0	17,19,21	0.63	0
2	NAG	c	2	2	14,14,15	0.21	0	17,19,21	0.62	0
5	NAG	d	1	5,1	14,14,15	0.36	0	17,19,21	0.75	1 (5%)
5	NAG	d	2	5	14,14,15	0.27	0	17,19,21	0.65	0
5	BMA	d	3	5	11,11,12	0.24	0	15,15,17	0.74	0
5	MAN	d	4	5	11,11,12	0.24	0	15,15,17	0.72	0
2	NAG	e	1	2,1	14,14,15	1.25	2 (14%)	17,19,21	2.63	7 (41%)
2	NAG	e	2	2	14,14,15	0.50	0	17,19,21	1.16	2 (11%)
2	NAG	f	1	2,1	14,14,15	0.19	0	17,19,21	0.72	1 (5%)
2	NAG	f	2	2	14,14,15	0.34	0	17,19,21	0.54	0
5	NAG	g	1	5,1	14,14,15	0.27	0	17,19,21	0.65	0
5	NAG	g	2	5	14,14,15	1.05	2 (14%)	17,19,21	1.97	5 (29%)
5	BMA	g	3	5	11,11,12	0.34	0	15,15,17	1.90	4 (26%)
5	MAN	g	4	5	11,11,12	0.41	0	15,15,17	0.92	1 (6%)
5	NAG	h	1	5,1	14,14,15	0.18	0	17,19,21	0.74	1 (5%)
5	NAG	h	2	5	14,14,15	0.24	0	17,19,21	0.58	0
5	BMA	h	3	5	11,11,12	0.25	0	15,15,17	0.52	0
5	MAN	h	4	5	11,11,12	0.28	0	15,15,17	0.53	0
2	NAG	i	1	2,1	14,14,15	0.37	0	17,19,21	0.52	0
2	NAG	i	2	2	14,14,15	0.32	0	17,19,21	0.58	0
6	NAG	j	1	6,1	14,14,15	0.31	0	17,19,21	0.88	0
6	NAG	j	2	6	14,14,15	0.65	0	17,19,21	1.89	2 (11%)
6	BMA	j	3	6	11,11,12	0.59	0	15,15,17	1.09	1 (6%)
6	FUC	j	4	6	10,10,11	0.79	0	14,14,16	0.73	0
7	NAG	k	1	7,1	14,14,15	0.32	0	17,19,21	0.59	0
7	NAG	k	2	7	14,14,15	0.21	0	17,19,21	0.57	0
7	BMA	k	3	7	11,11,12	0.57	0	15,15,17	2.36	2 (13%)
7	MAN	k	4	7	11,11,12	0.56	0	15,15,17	3.21	4 (26%)
7	MAN	k	5	7	11,11,12	0.26	0	15,15,17	0.51	0
8	NAG	l	1	8,1	14,14,15	2.09	7 (50%)	17,19,21	2.37	9 (52%)
8	NAG	l	2	8	14,14,15	0.28	0	17,19,21	0.94	1 (5%)
8	BMA	l	3	8	11,11,12	0.21	0	15,15,17	1.23	2 (13%)
8	MAN	l	4	8	11,11,12	0.26	0	15,15,17	0.94	1 (6%)
8	MAN	l	5	8	11,11,12	0.29	0	15,15,17	1.18	3 (20%)
8	MAN	l	6	8	11,11,12	0.31	0	15,15,17	0.84	0
8	MAN	l	7	8	11,11,12	0.96	1 (9%)	15,15,17	0.96	2 (13%)
2	NAG	m	1	2,1	14,14,15	0.18	0	17,19,21	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	m	2	2	14,14,15	0.26	0	17,19,21	0.52	0
2	NAG	n	1	2,1	14,14,15	0.30	0	17,19,21	0.56	0
2	NAG	n	2	2	14,14,15	0.29	0	17,19,21	0.81	0
3	NAG	o	1	3,1	14,14,15	0.45	0	17,19,21	1.99	6 (35%)
3	NAG	o	2	3	14,14,15	0.49	0	17,19,21	1.39	2 (11%)
3	BMA	o	3	3	11,11,12	0.43	0	15,15,17	1.36	2 (13%)
2	NAG	p	1	2,1	14,14,15	0.32	0	17,19,21	0.67	0
2	NAG	p	2	2	14,14,15	0.33	0	17,19,21	0.89	0
3	NAG	q	1	3,1	14,14,15	0.33	0	17,19,21	0.54	0
3	NAG	q	2	3	14,14,15	0.25	0	17,19,21	0.59	0
3	BMA	q	3	3	11,11,12	0.23	0	15,15,17	0.76	0
3	NAG	r	1	3,1	14,14,15	0.30	0	17,19,21	0.85	0
3	NAG	r	2	3	14,14,15	0.28	0	17,19,21	0.59	0
3	BMA	r	3	3	11,11,12	0.24	0	15,15,17	0.75	0
4	NAG	s	1	4,1	14,14,15	0.68	0	17,19,21	1.94	4 (23%)
4	NAG	s	2	4	14,14,15	0.49	0	17,19,21	1.68	4 (23%)
4	BMA	s	3	4	11,11,12	0.60	0	15,15,17	2.24	8 (53%)
4	MAN	s	4	4	11,11,12	0.68	0	15,15,17	0.96	0
4	MAN	s	5	4	11,11,12	0.60	0	15,15,17	1.38	2 (13%)
4	MAN	s	6	4	11,11,12	0.67	0	15,15,17	1.23	1 (6%)
4	MAN	s	7	4	11,11,12	0.67	0	15,15,17	0.94	0
5	NAG	t	1	5,1	14,14,15	0.54	0	17,19,21	1.27	3 (17%)
5	NAG	t	2	5	14,14,15	1.06	2 (14%)	17,19,21	1.85	3 (17%)
5	BMA	t	3	5	11,11,12	0.58	0	15,15,17	1.42	2 (13%)
5	MAN	t	4	5	11,11,12	0.37	0	15,15,17	1.23	2 (13%)
2	NAG	u	1	2,1	14,14,15	0.24	0	17,19,21	0.63	0
2	NAG	u	2	2	14,14,15	0.21	0	17,19,21	0.62	0
5	NAG	v	1	5,1	14,14,15	0.37	0	17,19,21	0.75	1 (5%)
5	NAG	v	2	5	14,14,15	0.27	0	17,19,21	0.65	0
5	BMA	v	3	5	11,11,12	0.24	0	15,15,17	0.73	0
5	MAN	v	4	5	11,11,12	0.25	0	15,15,17	0.73	0
2	NAG	w	1	2,1	14,14,15	1.25	2 (14%)	17,19,21	2.63	7 (41%)
2	NAG	w	2	2	14,14,15	0.50	0	17,19,21	1.16	2 (11%)
2	NAG	x	1	2,1	14,14,15	0.18	0	17,19,21	0.72	1 (5%)
2	NAG	x	2	2	14,14,15	0.35	0	17,19,21	0.54	0
5	NAG	y	1	5,1	14,14,15	0.27	0	17,19,21	0.65	0
5	NAG	y	2	5	14,14,15	1.05	1 (7%)	17,19,21	1.97	4 (23%)
5	BMA	y	3	5	11,11,12	0.34	0	15,15,17	1.90	4 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MAN	y	4	5	11,11,12	0.42	0	15,15,17	0.91	1 (6%)
5	NAG	z	1	5,1	14,14,15	0.18	0	17,19,21	0.74	1 (5%)
5	NAG	z	2	5	14,14,15	0.24	0	17,19,21	0.57	0
5	BMA	z	3	5	11,11,12	0.25	0	15,15,17	0.52	0
5	MAN	z	4	5	11,11,12	0.28	0	15,15,17	0.52	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	0	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	0	2	2	-	2/6/23/26	0/1/1/1
6	NAG	1	1	6,1	-	4/6/23/26	0/1/1/1
6	NAG	1	2	6	-	0/6/23/26	0/1/1/1
6	BMA	1	3	6	-	0/2/19/22	0/1/1/1
6	FUC	1	4	6	-	-	0/1/1/1
7	NAG	2	1	7,1	-	3/6/23/26	0/1/1/1
7	NAG	2	2	7	-	0/6/23/26	0/1/1/1
7	BMA	2	3	7	-	0/2/19/22	0/1/1/1
7	MAN	2	4	7	-	0/2/19/22	0/1/1/1
7	MAN	2	5	7	-	0/2/19/22	0/1/1/1
8	NAG	3	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	3	2	8	-	3/6/23/26	0/1/1/1
8	BMA	3	3	8	-	1/2/19/22	0/1/1/1
8	MAN	3	4	8	-	1/2/19/22	0/1/1/1
8	MAN	3	5	8	-	2/2/19/22	0/1/1/1
8	MAN	3	6	8	-	0/2/19/22	0/1/1/1
8	MAN	3	7	8	-	2/2/19/22	0/1/1/1
2	NAG	4	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	4	2	2	-	0/6/23/26	0/1/1/1
2	NAG	D	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	1/6/23/26	0/1/1/1
3	NAG	E	1	3,1	-	5/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
2	NAG	F	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	G	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	3/6/23/26	0/1/1/1
3	BMA	G	3	3	-	1/2/19/22	0/1/1/1
3	NAG	H	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
3	BMA	H	3	3	-	0/2/19/22	0/1/1/1
4	NAG	I	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	I	2	4	-	3/6/23/26	0/1/1/1
4	BMA	I	3	4	-	2/2/19/22	0/1/1/1
4	MAN	I	4	4	-	0/2/19/22	0/1/1/1
4	MAN	I	5	4	-	1/2/19/22	0/1/1/1
4	MAN	I	6	4	-	2/2/19/22	0/1/1/1
4	MAN	I	7	4	-	0/2/19/22	0/1/1/1
5	NAG	J	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	J	2	5	-	2/6/23/26	0/1/1/1
5	BMA	J	3	5	-	1/2/19/22	0/1/1/1
5	MAN	J	4	5	-	0/2/19/22	0/1/1/1
2	NAG	K	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	K	2	2	-	2/6/23/26	0/1/1/1
5	NAG	L	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	L	2	5	-	1/6/23/26	0/1/1/1
5	BMA	L	3	5	-	0/2/19/22	0/1/1/1
5	MAN	L	4	5	-	1/2/19/22	0/1/1/1
2	NAG	M	1	2,1	-	3/6/23/26	0/1/1/1
2	NAG	M	2	2	-	3/6/23/26	0/1/1/1
2	NAG	N	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	N	2	2	-	2/6/23/26	0/1/1/1
5	NAG	O	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	O	2	5	-	1/6/23/26	0/1/1/1
5	BMA	O	3	5	-	2/2/19/22	0/1/1/1
5	MAN	O	4	5	-	0/2/19/22	0/1/1/1
5	NAG	P	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	P	2	5	-	2/6/23/26	0/1/1/1
5	BMA	P	3	5	-	0/2/19/22	0/1/1/1
5	MAN	P	4	5	-	0/2/19/22	0/1/1/1
2	NAG	Q	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	Q	2	2	-	2/6/23/26	0/1/1/1

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	b	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	b	2	5	-	2/6/23/26	0/1/1/1
5	BMA	b	3	5	-	1/2/19/22	0/1/1/1
5	MAN	b	4	5	-	0/2/19/22	0/1/1/1
2	NAG	c	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	c	2	2	-	2/6/23/26	0/1/1/1
5	NAG	d	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	d	2	5	-	1/6/23/26	0/1/1/1
5	BMA	d	3	5	-	0/2/19/22	0/1/1/1
5	MAN	d	4	5	-	1/2/19/22	0/1/1/1
2	NAG	e	1	2,1	-	3/6/23/26	0/1/1/1
2	NAG	e	2	2	-	3/6/23/26	0/1/1/1
2	NAG	f	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	f	2	2	-	2/6/23/26	0/1/1/1
5	NAG	g	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	g	2	5	-	1/6/23/26	0/1/1/1
5	BMA	g	3	5	-	2/2/19/22	0/1/1/1
5	MAN	g	4	5	-	0/2/19/22	0/1/1/1
5	NAG	h	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	h	2	5	-	2/6/23/26	0/1/1/1
5	BMA	h	3	5	-	0/2/19/22	0/1/1/1
5	MAN	h	4	5	-	0/2/19/22	0/1/1/1
2	NAG	i	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	i	2	2	-	2/6/23/26	0/1/1/1
6	NAG	j	1	6,1	-	4/6/23/26	0/1/1/1
6	NAG	j	2	6	-	0/6/23/26	0/1/1/1
6	BMA	j	3	6	-	0/2/19/22	0/1/1/1
6	FUC	j	4	6	-	-	0/1/1/1
7	NAG	k	1	7,1	-	3/6/23/26	0/1/1/1
7	NAG	k	2	7	-	0/6/23/26	0/1/1/1
7	BMA	k	3	7	-	0/2/19/22	0/1/1/1
7	MAN	k	4	7	-	0/2/19/22	0/1/1/1
7	MAN	k	5	7	-	0/2/19/22	0/1/1/1
8	NAG	l	1	8,1	-	2/6/23/26	0/1/1/1
8	NAG	l	2	8	-	3/6/23/26	0/1/1/1
8	BMA	l	3	8	-	1/2/19/22	0/1/1/1
8	MAN	l	4	8	-	1/2/19/22	0/1/1/1
8	MAN	l	5	8	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	MAN	l	6	8	-	0/2/19/22	0/1/1/1
8	MAN	l	7	8	-	2/2/19/22	0/1/1/1
2	NAG	m	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	m	2	2	-	0/6/23/26	0/1/1/1
2	NAG	n	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	n	2	2	-	1/6/23/26	0/1/1/1
3	NAG	o	1	3,1	-	5/6/23/26	0/1/1/1
3	NAG	o	2	3	-	2/6/23/26	0/1/1/1
3	BMA	o	3	3	-	0/2/19/22	0/1/1/1
2	NAG	p	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	p	2	2	-	2/6/23/26	0/1/1/1
3	NAG	q	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	q	2	3	-	3/6/23/26	0/1/1/1
3	BMA	q	3	3	-	1/2/19/22	0/1/1/1
3	NAG	r	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	r	2	3	-	0/6/23/26	0/1/1/1
3	BMA	r	3	3	-	0/2/19/22	0/1/1/1
4	NAG	s	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	s	2	4	-	3/6/23/26	0/1/1/1
4	BMA	s	3	4	-	2/2/19/22	0/1/1/1
4	MAN	s	4	4	-	0/2/19/22	0/1/1/1
4	MAN	s	5	4	-	1/2/19/22	0/1/1/1
4	MAN	s	6	4	-	2/2/19/22	0/1/1/1
4	MAN	s	7	4	-	0/2/19/22	0/1/1/1
5	NAG	t	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	t	2	5	-	2/6/23/26	0/1/1/1
5	BMA	t	3	5	-	1/2/19/22	0/1/1/1
5	MAN	t	4	5	-	0/2/19/22	0/1/1/1
2	NAG	u	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	u	2	2	-	2/6/23/26	0/1/1/1
5	NAG	v	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	v	2	5	-	1/6/23/26	0/1/1/1
5	BMA	v	3	5	-	0/2/19/22	0/1/1/1
5	MAN	v	4	5	-	1/2/19/22	0/1/1/1
2	NAG	w	1	2,1	-	3/6/23/26	0/1/1/1
2	NAG	w	2	2	-	3/6/23/26	0/1/1/1
2	NAG	x	1	2,1	-	4/6/23/26	0/1/1/1

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

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	3	1	NAG	O5-C1	-4.43	1.36	1.43
8	l	1	NAG	O5-C1	-4.42	1.36	1.43
8	T	1	NAG	O5-C1	-4.42	1.36	1.43
8	T	1	NAG	C1-C2	-3.27	1.47	1.52
8	l	1	NAG	C1-C2	-3.24	1.47	1.52
8	3	1	NAG	C1-C2	-3.21	1.48	1.52
8	T	1	NAG	C2-N2	-2.64	1.41	1.46
8	l	1	NAG	C2-N2	-2.64	1.41	1.46
8	3	1	NAG	C2-N2	-2.64	1.41	1.46
2	w	1	NAG	C3-C2	-2.52	1.47	1.52
2	e	1	NAG	C3-C2	-2.50	1.47	1.52
2	M	1	NAG	C3-C2	-2.50	1.47	1.52
8	3	1	NAG	C4-C3	-2.45	1.46	1.52
8	l	1	NAG	C4-C3	-2.44	1.46	1.52
8	T	1	NAG	C4-C3	-2.44	1.46	1.52
8	l	1	NAG	C3-C2	-2.35	1.47	1.52
5	y	2	NAG	C2-N2	-2.35	1.42	1.46
8	3	7	MAN	C1-C2	2.34	1.57	1.52
8	3	1	NAG	C3-C2	-2.33	1.47	1.52
8	T	1	NAG	C3-C2	-2.33	1.47	1.52
8	T	7	MAN	C1-C2	2.33	1.57	1.52
8	l	7	MAN	C1-C2	2.32	1.57	1.52
5	g	2	NAG	C2-N2	-2.32	1.42	1.46
5	O	2	NAG	C2-N2	-2.30	1.42	1.46
2	M	1	NAG	O5-C1	-2.22	1.40	1.43
8	T	1	NAG	O5-C5	-2.22	1.39	1.43
2	e	1	NAG	O5-C1	-2.21	1.40	1.43
2	w	1	NAG	O5-C1	-2.21	1.40	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	b	2	NAG	C1-C2	-2.20	1.49	1.52
5	t	2	NAG	C1-C2	-2.19	1.49	1.52
5	J	2	NAG	C1-C2	-2.19	1.49	1.52
8	3	1	NAG	O5-C5	-2.19	1.39	1.43
8	l	1	NAG	O5-C5	-2.18	1.39	1.43
8	3	1	NAG	C4-C5	-2.07	1.48	1.53
8	l	1	NAG	C4-C5	-2.07	1.48	1.53
8	T	1	NAG	C4-C5	-2.06	1.48	1.53
5	g	2	NAG	C1-C2	-2.02	1.49	1.52
5	t	2	NAG	C2-N2	-2.01	1.42	1.46

All (264) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	S	3	BMA	O2-C2-C3	8.22	127.17	110.15
7	2	3	BMA	O2-C2-C3	8.22	127.17	110.15
7	k	3	BMA	O2-C2-C3	8.21	127.16	110.15
7	2	4	MAN	O2-C2-C3	7.82	126.35	110.15
7	k	4	MAN	O2-C2-C3	7.82	126.34	110.15
7	S	4	MAN	O2-C2-C3	7.80	126.31	110.15
6	R	2	NAG	O4-C4-C3	-6.76	94.43	110.38
6	j	2	NAG	O4-C4-C3	-6.76	94.45	110.38
6	1	2	NAG	O4-C4-C3	-6.76	94.45	110.38
5	b	2	NAG	C2-N2-C7	-6.09	114.73	122.90
5	J	2	NAG	C2-N2-C7	-6.08	114.75	122.90
5	t	2	NAG	C2-N2-C7	-6.07	114.76	122.90
7	2	4	MAN	O3-C3-C4	-5.88	96.51	110.38
7	S	4	MAN	O3-C3-C4	-5.88	96.51	110.38
7	k	4	MAN	O3-C3-C4	-5.87	96.53	110.38
7	2	4	MAN	O4-C4-C5	-5.81	95.02	109.32
7	k	4	MAN	O4-C4-C5	-5.81	95.02	109.32
2	M	1	NAG	O3-C3-C2	-5.80	97.34	109.40
2	e	1	NAG	O3-C3-C2	-5.80	97.34	109.40
7	S	4	MAN	O4-C4-C5	-5.80	95.04	109.32
2	w	1	NAG	O3-C3-C2	-5.80	97.36	109.40
4	s	3	BMA	O3-C3-C4	4.91	121.96	110.38
4	a	3	BMA	O3-C3-C4	4.91	121.95	110.38
4	I	3	BMA	O3-C3-C4	4.90	121.93	110.38
3	W	1	NAG	C4-C3-C2	-4.50	104.42	111.02
3	E	1	NAG	C4-C3-C2	-4.49	104.44	111.02
3	o	1	NAG	C4-C3-C2	-4.48	104.45	111.02
8	T	1	NAG	C6-C5-C4	-4.45	102.10	113.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	3	1	NAG	C6-C5-C4	-4.44	102.11	113.02
8	l	1	NAG	C6-C5-C4	-4.44	102.12	113.02
7	2	4	MAN	O5-C5-C6	-4.31	99.27	107.66
7	k	4	MAN	O5-C5-C6	-4.31	99.28	107.66
4	a	1	NAG	O5-C1-C2	-4.30	104.64	111.29
7	S	4	MAN	O5-C5-C6	-4.30	99.30	107.66
4	s	1	NAG	O5-C1-C2	-4.29	104.66	111.29
4	I	1	NAG	O5-C1-C2	-4.29	104.66	111.29
4	I	1	NAG	C1-O5-C5	4.27	117.92	112.19
4	a	1	NAG	C1-O5-C5	4.26	117.90	112.19
4	s	1	NAG	C1-O5-C5	4.25	117.89	112.19
2	e	1	NAG	O4-C4-C5	-4.21	98.96	109.32
2	M	1	NAG	O4-C4-C5	-4.21	98.96	109.32
2	w	1	NAG	O4-C4-C5	-4.20	98.97	109.32
5	O	2	NAG	O5-C5-C6	4.15	115.75	107.66
5	g	2	NAG	O5-C5-C6	4.14	115.72	107.66
5	y	2	NAG	O5-C5-C6	4.14	115.72	107.66
5	O	2	NAG	C2-N2-C7	-4.04	117.48	122.90
2	e	1	NAG	C1-O5-C5	4.02	117.58	112.19
2	w	1	NAG	C1-O5-C5	4.01	117.56	112.19
2	M	1	NAG	O3-C3-C4	-4.00	100.95	110.38
5	y	2	NAG	C2-N2-C7	-4.00	117.54	122.90
2	M	1	NAG	C1-O5-C5	4.00	117.54	112.19
2	w	1	NAG	O3-C3-C4	-4.00	100.95	110.38
5	g	2	NAG	C2-N2-C7	-3.99	117.55	122.90
2	e	1	NAG	O3-C3-C4	-3.99	100.97	110.38
8	3	1	NAG	C1-C2-N2	-3.97	104.17	110.43
8	T	1	NAG	C1-C2-N2	-3.97	104.18	110.43
8	l	1	NAG	C1-C2-N2	-3.97	104.18	110.43
4	s	2	NAG	C2-N2-C7	-3.89	117.69	122.90
4	I	2	NAG	C2-N2-C7	-3.88	117.70	122.90
4	a	2	NAG	C2-N2-C7	-3.86	117.73	122.90
5	g	3	BMA	O5-C5-C6	3.66	114.79	107.66
5	O	3	BMA	O5-C5-C6	3.66	114.78	107.66
5	y	3	BMA	O5-C5-C6	3.65	114.77	107.66
3	o	1	NAG	O4-C4-C3	3.57	118.79	110.38
3	W	1	NAG	O4-C4-C3	3.56	118.77	110.38
3	E	1	NAG	O4-C4-C3	3.56	118.77	110.38
8	3	1	NAG	O3-C3-C2	-3.38	102.37	109.40
8	T	1	NAG	O3-C3-C2	-3.37	102.40	109.40
8	l	1	NAG	O3-C3-C2	-3.37	102.41	109.40
5	y	3	BMA	C1-O5-C5	-3.34	107.71	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	O	3	BMA	C1-O5-C5	-3.34	107.71	112.19
5	g	3	BMA	C1-O5-C5	-3.33	107.72	112.19
4	I	1	NAG	C2-N2-C7	-3.23	118.57	122.90
4	s	1	NAG	C2-N2-C7	-3.23	118.58	122.90
4	a	1	NAG	C2-N2-C7	-3.21	118.60	122.90
3	o	3	BMA	C2-C3-C4	-3.18	105.28	110.86
3	E	3	BMA	C2-C3-C4	-3.14	105.33	110.86
3	W	3	BMA	C2-C3-C4	-3.14	105.34	110.86
3	o	2	NAG	C1-O5-C5	3.11	116.36	112.19
4	s	2	NAG	O4-C4-C5	-3.10	101.69	109.32
4	I	2	NAG	O4-C4-C5	-3.09	101.70	109.32
4	a	2	NAG	O4-C4-C5	-3.09	101.72	109.32
3	E	2	NAG	C1-O5-C5	3.08	116.32	112.19
3	W	2	NAG	C1-O5-C5	3.06	116.28	112.19
5	O	2	NAG	C1-C2-N2	-3.03	105.67	110.43
5	g	2	NAG	C1-C2-N2	-3.02	105.67	110.43
5	y	2	NAG	C1-C2-N2	-3.00	105.70	110.43
4	I	5	MAN	C2-C3-C4	-2.94	105.69	110.86
4	s	5	MAN	C2-C3-C4	-2.94	105.69	110.86
8	l	1	NAG	O4-C4-C5	-2.94	102.09	109.32
8	T	1	NAG	O4-C4-C5	-2.94	102.09	109.32
4	a	5	MAN	C2-C3-C4	-2.94	105.69	110.86
8	3	1	NAG	O4-C4-C5	-2.94	102.09	109.32
4	a	3	BMA	C2-C3-C4	-2.89	105.79	110.86
4	I	3	BMA	C2-C3-C4	-2.88	105.80	110.86
4	s	3	BMA	C2-C3-C4	-2.88	105.80	110.86
3	E	1	NAG	O5-C1-C2	-2.87	106.86	111.29
3	E	1	NAG	O5-C5-C6	2.85	113.22	107.66
3	W	1	NAG	C2-N2-C7	-2.85	119.08	122.90
3	W	1	NAG	O5-C1-C2	-2.85	106.88	111.29
3	o	1	NAG	O5-C5-C6	2.85	113.21	107.66
3	o	1	NAG	O5-C1-C2	-2.85	106.89	111.29
3	W	1	NAG	O5-C5-C6	2.84	113.20	107.66
5	b	1	NAG	C2-N2-C7	-2.84	119.09	122.90
3	o	1	NAG	C2-N2-C7	-2.84	119.09	122.90
3	E	1	NAG	C2-N2-C7	-2.83	119.10	122.90
5	t	1	NAG	C2-N2-C7	-2.83	119.10	122.90
5	g	3	BMA	C3-C4-C5	-2.82	105.11	110.23
5	J	1	NAG	C2-N2-C7	-2.81	119.13	122.90
5	y	3	BMA	C3-C4-C5	-2.81	105.13	110.23
8	l	5	MAN	C1-C2-C3	2.81	113.74	109.64
8	3	5	MAN	C1-C2-C3	2.81	113.74	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	T	5	MAN	C1-C2-C3	2.81	113.73	109.64
2	M	1	NAG	C6-C5-C4	-2.80	106.14	113.02
2	w	1	NAG	C6-C5-C4	-2.80	106.14	113.02
2	e	1	NAG	C6-C5-C4	-2.80	106.15	113.02
5	O	3	BMA	C3-C4-C5	-2.80	105.16	110.23
6	l	2	NAG	O4-C4-C5	-2.73	102.61	109.32
6	j	2	NAG	O4-C4-C5	-2.72	102.62	109.32
6	R	2	NAG	O4-C4-C5	-2.72	102.62	109.32
4	s	3	BMA	C1-O5-C5	-2.69	108.58	112.19
4	a	3	BMA	C1-O5-C5	-2.69	108.58	112.19
4	I	3	BMA	C3-C4-C5	-2.68	105.37	110.23
4	I	3	BMA	C1-O5-C5	-2.68	108.60	112.19
4	a	3	BMA	C3-C4-C5	-2.67	105.39	110.23
4	s	3	BMA	C3-C4-C5	-2.67	105.39	110.23
4	I	3	BMA	O5-C5-C6	2.63	112.79	107.66
4	a	3	BMA	O5-C5-C6	2.63	112.78	107.66
4	s	3	BMA	O5-C5-C6	2.62	112.77	107.66
5	J	3	BMA	C6-C5-C4	-2.62	106.58	113.02
5	b	3	BMA	C6-C5-C4	-2.61	106.61	113.02
5	t	3	BMA	C6-C5-C4	-2.61	106.61	113.02
4	s	2	NAG	O5-C1-C2	-2.60	107.26	111.29
4	I	2	NAG	O5-C1-C2	-2.60	107.26	111.29
4	a	2	NAG	O5-C1-C2	-2.60	107.27	111.29
8	3	1	NAG	C2-N2-C7	-2.57	119.45	122.90
8	T	1	NAG	C2-N2-C7	-2.56	119.47	122.90
8	l	1	NAG	C2-N2-C7	-2.56	119.47	122.90
5	J	4	MAN	C2-C3-C4	-2.52	106.42	110.86
5	t	4	MAN	C2-C3-C4	-2.52	106.44	110.86
5	b	4	MAN	C2-C3-C4	-2.51	106.45	110.86
5	g	3	BMA	C2-C3-C4	-2.50	106.46	110.86
5	g	4	MAN	C2-C3-C4	-2.50	106.47	110.86
5	y	3	BMA	C2-C3-C4	-2.49	106.48	110.86
5	O	3	BMA	C2-C3-C4	-2.49	106.48	110.86
5	O	4	MAN	C2-C3-C4	-2.49	106.49	110.86
2	M	1	NAG	O5-C5-C6	-2.48	102.84	107.66
2	e	1	NAG	O5-C5-C6	-2.47	102.86	107.66
2	w	1	NAG	O5-C5-C6	-2.47	102.86	107.66
4	a	5	MAN	O2-C2-C3	-2.46	105.06	110.15
5	y	4	MAN	C2-C3-C4	-2.46	106.54	110.86
4	I	5	MAN	O2-C2-C3	-2.45	105.08	110.15
4	s	5	MAN	O2-C2-C3	-2.45	105.08	110.15
8	l	1	NAG	O5-C1-C2	-2.44	107.51	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	3	1	NAG	O5-C1-C2	-2.43	107.53	111.29
8	l	1	NAG	O5-C5-C6	-2.43	102.94	107.66
8	T	1	NAG	O5-C1-C2	-2.42	107.54	111.29
8	3	1	NAG	O5-C5-C6	-2.42	102.95	107.66
8	T	1	NAG	O5-C5-C6	-2.42	102.96	107.66
7	2	3	BMA	C2-C3-C4	2.40	115.08	110.86
7	k	3	BMA	C2-C3-C4	2.38	115.05	110.86
7	S	3	BMA	C2-C3-C4	2.38	115.04	110.86
6	j	3	BMA	C2-C3-C4	2.36	115.01	110.86
6	R	3	BMA	C2-C3-C4	2.35	115.00	110.86
6	l	3	BMA	C2-C3-C4	2.35	115.00	110.86
5	y	2	NAG	C4-C3-C2	-2.34	107.59	111.02
5	O	2	NAG	C4-C3-C2	-2.34	107.59	111.02
3	W	3	BMA	O5-C1-C2	-2.34	105.22	110.79
5	g	2	NAG	C4-C3-C2	-2.34	107.59	111.02
8	T	2	NAG	O5-C1-C2	-2.33	107.68	111.29
4	s	3	BMA	O4-C4-C3	2.33	115.87	110.38
3	E	3	BMA	O5-C1-C2	-2.33	105.23	110.79
5	t	3	BMA	C2-C3-C4	-2.33	106.77	110.86
4	a	3	BMA	O4-C4-C3	2.33	115.86	110.38
8	l	2	NAG	O5-C1-C2	-2.33	107.69	111.29
3	o	3	BMA	O5-C1-C2	-2.33	105.24	110.79
8	T	1	NAG	C4-C3-C2	-2.32	107.61	111.02
8	3	1	NAG	C4-C3-C2	-2.32	107.61	111.02
5	b	3	BMA	C2-C3-C4	-2.32	106.78	110.86
4	I	3	BMA	O4-C4-C3	2.32	115.84	110.38
8	l	1	NAG	C4-C3-C2	-2.31	107.63	111.02
5	J	3	BMA	C2-C3-C4	-2.31	106.80	110.86
8	3	2	NAG	O5-C1-C2	-2.29	107.75	111.29
4	a	1	NAG	C3-C4-C5	-2.28	106.10	110.23
4	s	1	NAG	C3-C4-C5	-2.28	106.10	110.23
4	I	1	NAG	C3-C4-C5	-2.27	106.11	110.23
8	3	5	MAN	O5-C1-C2	2.25	116.17	110.79
8	l	5	MAN	O5-C1-C2	2.25	116.16	110.79
8	T	5	MAN	O5-C1-C2	2.25	116.14	110.79
8	T	5	MAN	C1-O5-C5	2.24	115.18	112.19
3	o	2	NAG	O4-C4-C3	-2.24	105.11	110.38
3	W	2	NAG	O4-C4-C3	-2.23	105.12	110.38
3	E	2	NAG	O4-C4-C3	-2.22	105.13	110.38
3	E	1	NAG	C1-O5-C5	2.22	115.17	112.19
4	s	3	BMA	O4-C4-C5	-2.22	103.86	109.32
2	w	2	NAG	C4-C3-C2	-2.22	107.77	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	3	5	MAN	C1-O5-C5	2.22	115.16	112.19
4	I	3	BMA	O4-C4-C5	-2.22	103.86	109.32
2	M	2	NAG	C4-C3-C2	-2.22	107.77	111.02
5	t	1	NAG	C6-C5-C4	-2.21	107.58	113.02
5	t	2	NAG	O4-C4-C3	-2.21	105.16	110.38
8	l	5	MAN	C1-O5-C5	2.21	115.15	112.19
2	e	2	NAG	C4-C3-C2	-2.21	107.78	111.02
5	b	2	NAG	O4-C4-C3	-2.21	105.17	110.38
5	J	1	NAG	C6-C5-C4	-2.21	107.59	113.02
5	b	1	NAG	C6-C5-C4	-2.21	107.59	113.02
4	a	3	BMA	O4-C4-C5	-2.21	103.89	109.32
5	J	2	NAG	O4-C4-C3	-2.20	105.18	110.38
8	l	1	NAG	C3-C4-C5	-2.20	106.24	110.23
8	3	1	NAG	C3-C4-C5	-2.20	106.24	110.23
8	T	1	NAG	C3-C4-C5	-2.20	106.24	110.23
8	l	3	BMA	C2-C3-C4	-2.20	106.99	110.86
3	W	1	NAG	C1-O5-C5	2.20	115.13	112.19
8	T	3	BMA	C2-C3-C4	-2.20	107.00	110.86
8	3	3	BMA	C2-C3-C4	-2.19	107.01	110.86
3	o	1	NAG	C1-O5-C5	2.18	115.11	112.19
8	T	7	MAN	C1-O5-C5	2.18	115.11	112.19
8	l	7	MAN	C1-O5-C5	2.18	115.10	112.19
8	3	7	MAN	C1-O5-C5	2.17	115.10	112.19
5	J	2	NAG	O5-C1-C2	-2.17	107.93	111.29
5	t	1	NAG	C1-C2-N2	2.16	113.84	110.43
5	b	2	NAG	O5-C1-C2	-2.16	107.94	111.29
2	e	1	NAG	O5-C1-C2	-2.16	107.94	111.29
2	w	1	NAG	O5-C1-C2	-2.16	107.95	111.29
2	M	2	NAG	O5-C1-C2	-2.16	107.95	111.29
5	t	2	NAG	O5-C1-C2	-2.15	107.96	111.29
5	J	1	NAG	C1-C2-N2	2.15	113.82	110.43
2	x	1	NAG	C1-O5-C5	2.15	115.07	112.19
5	b	1	NAG	C1-C2-N2	2.15	113.82	110.43
2	w	2	NAG	O5-C1-C2	-2.15	107.97	111.29
2	M	1	NAG	O5-C1-C2	-2.15	107.97	111.29
2	e	2	NAG	O5-C1-C2	-2.15	107.97	111.29
2	f	1	NAG	C1-O5-C5	2.14	115.05	112.19
2	N	1	NAG	C1-O5-C5	2.13	115.04	112.19
8	l	3	BMA	O5-C5-C6	2.12	111.80	107.66
5	h	1	NAG	C1-O5-C5	2.12	115.03	112.19
5	P	1	NAG	C1-O5-C5	2.12	115.02	112.19
5	z	1	NAG	C1-O5-C5	2.11	115.01	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	T	3	BMA	O5-C5-C6	2.11	111.76	107.66
5	b	4	MAN	O5-C1-C2	-2.10	105.77	110.79
8	3	3	BMA	O5-C5-C6	2.10	111.75	107.66
4	a	6	MAN	C1-O5-C5	2.09	114.99	112.19
8	3	7	MAN	O2-C2-C3	-2.09	105.83	110.15
4	I	6	MAN	C1-O5-C5	2.09	114.98	112.19
8	l	7	MAN	O2-C2-C3	-2.08	105.83	110.15
5	J	4	MAN	O5-C1-C2	-2.08	105.82	110.79
4	s	6	MAN	C1-O5-C5	2.08	114.97	112.19
8	T	7	MAN	O2-C2-C3	-2.08	105.84	110.15
5	t	4	MAN	O5-C1-C2	-2.08	105.84	110.79
5	L	1	NAG	C1-O5-C5	2.07	114.96	112.19
5	d	1	NAG	C1-O5-C5	2.05	114.94	112.19
5	v	1	NAG	C1-O5-C5	2.05	114.93	112.19
8	3	4	MAN	C2-C3-C4	-2.04	107.28	110.86
4	I	2	NAG	C3-C4-C5	-2.03	106.54	110.23
8	T	4	MAN	C2-C3-C4	-2.03	107.28	110.86
4	s	2	NAG	C3-C4-C5	-2.03	106.55	110.23
4	I	3	BMA	O2-C2-C1	-2.03	104.58	109.22
4	a	2	NAG	C3-C4-C5	-2.02	106.56	110.23
8	l	4	MAN	C2-C3-C4	-2.02	107.31	110.86
4	s	3	BMA	O2-C2-C1	-2.02	104.61	109.22
2	Q	2	NAG	C1-O5-C5	2.02	114.89	112.19
2	0	2	NAG	C1-O5-C5	2.01	114.88	112.19
4	a	3	BMA	O2-C2-C1	-2.01	104.62	109.22
5	g	2	NAG	C3-C4-C5	-2.01	106.59	110.23

There are no chirality outliers.

All (288) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	M	1	NAG	C3-C2-N2-C7
2	M	2	NAG	C8-C7-N2-C2
2	M	2	NAG	O7-C7-N2-C2
2	e	1	NAG	C3-C2-N2-C7
2	e	2	NAG	C8-C7-N2-C2
2	e	2	NAG	O7-C7-N2-C2
2	w	1	NAG	C3-C2-N2-C7
2	w	2	NAG	C8-C7-N2-C2
2	w	2	NAG	O7-C7-N2-C2
3	E	1	NAG	C8-C7-N2-C2
3	E	1	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	G	2	NAG	C1-C2-N2-C7
3	G	2	NAG	C8-C7-N2-C2
3	G	2	NAG	O7-C7-N2-C2
3	H	1	NAG	C8-C7-N2-C2
3	H	1	NAG	O7-C7-N2-C2
3	W	1	NAG	C8-C7-N2-C2
3	W	1	NAG	O7-C7-N2-C2
3	Y	2	NAG	C1-C2-N2-C7
3	Y	2	NAG	C8-C7-N2-C2
3	Y	2	NAG	O7-C7-N2-C2
3	Z	1	NAG	C8-C7-N2-C2
3	Z	1	NAG	O7-C7-N2-C2
3	o	1	NAG	C8-C7-N2-C2
3	o	1	NAG	O7-C7-N2-C2
3	q	2	NAG	C1-C2-N2-C7
3	q	2	NAG	C8-C7-N2-C2
3	q	2	NAG	O7-C7-N2-C2
3	r	1	NAG	C8-C7-N2-C2
3	r	1	NAG	O7-C7-N2-C2
4	I	1	NAG	C8-C7-N2-C2
4	I	1	NAG	O7-C7-N2-C2
4	I	2	NAG	C3-C2-N2-C7
4	I	2	NAG	C8-C7-N2-C2
4	I	2	NAG	O7-C7-N2-C2
4	a	1	NAG	C8-C7-N2-C2
4	a	1	NAG	O7-C7-N2-C2
4	a	2	NAG	C3-C2-N2-C7
4	a	2	NAG	C8-C7-N2-C2
4	a	2	NAG	O7-C7-N2-C2
4	s	1	NAG	C8-C7-N2-C2
4	s	1	NAG	O7-C7-N2-C2
4	s	2	NAG	C3-C2-N2-C7
4	s	2	NAG	C8-C7-N2-C2
4	s	2	NAG	O7-C7-N2-C2
5	J	2	NAG	O7-C7-N2-C2
5	b	2	NAG	O7-C7-N2-C2
5	t	2	NAG	O7-C7-N2-C2
8	T	1	NAG	C1-C2-N2-C7
8	l	1	NAG	C1-C2-N2-C7
8	3	1	NAG	C1-C2-N2-C7
2	F	1	NAG	C8-C7-N2-C2
2	F	1	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	X	1	NAG	C8-C7-N2-C2
2	X	1	NAG	O7-C7-N2-C2
2	p	1	NAG	C8-C7-N2-C2
2	p	1	NAG	O7-C7-N2-C2
5	J	2	NAG	C8-C7-N2-C2
5	b	2	NAG	C8-C7-N2-C2
5	t	2	NAG	C8-C7-N2-C2
3	E	1	NAG	O5-C5-C6-O6
3	W	1	NAG	O5-C5-C6-O6
3	o	1	NAG	O5-C5-C6-O6
2	F	2	NAG	O7-C7-N2-C2
2	X	2	NAG	O7-C7-N2-C2
2	p	2	NAG	O7-C7-N2-C2
8	T	5	MAN	O5-C5-C6-O6
8	l	5	MAN	O5-C5-C6-O6
8	3	5	MAN	O5-C5-C6-O6
6	R	1	NAG	O5-C5-C6-O6
6	j	1	NAG	O5-C5-C6-O6
6	l	1	NAG	O5-C5-C6-O6
5	O	3	BMA	O5-C5-C6-O6
5	g	3	BMA	O5-C5-C6-O6
5	y	3	BMA	O5-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
2	c	2	NAG	O5-C5-C6-O6
2	u	2	NAG	O5-C5-C6-O6
4	I	6	MAN	O5-C5-C6-O6
4	a	6	MAN	O5-C5-C6-O6
4	s	6	MAN	O5-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
3	W	1	NAG	C4-C5-C6-O6
3	o	1	NAG	C4-C5-C6-O6
8	T	5	MAN	C4-C5-C6-O6
8	l	5	MAN	C4-C5-C6-O6
8	3	5	MAN	C4-C5-C6-O6
2	F	2	NAG	C8-C7-N2-C2
2	X	2	NAG	C8-C7-N2-C2
2	p	2	NAG	C8-C7-N2-C2
3	E	2	NAG	C8-C7-N2-C2
3	E	2	NAG	O7-C7-N2-C2
3	W	2	NAG	C8-C7-N2-C2
3	W	2	NAG	O7-C7-N2-C2
3	o	2	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	o	2	NAG	O7-C7-N2-C2
4	I	3	BMA	C4-C5-C6-O6
4	a	3	BMA	C4-C5-C6-O6
4	s	3	BMA	C4-C5-C6-O6
2	K	2	NAG	C4-C5-C6-O6
2	c	2	NAG	C4-C5-C6-O6
2	u	2	NAG	C4-C5-C6-O6
4	I	6	MAN	C4-C5-C6-O6
4	a	6	MAN	C4-C5-C6-O6
4	s	6	MAN	C4-C5-C6-O6
5	J	1	NAG	C8-C7-N2-C2
5	b	1	NAG	C8-C7-N2-C2
5	t	1	NAG	C8-C7-N2-C2
6	R	1	NAG	C4-C5-C6-O6
6	j	1	NAG	C4-C5-C6-O6
6	l	1	NAG	C4-C5-C6-O6
2	N	2	NAG	O5-C5-C6-O6
2	f	2	NAG	O5-C5-C6-O6
2	x	2	NAG	O5-C5-C6-O6
5	O	3	BMA	C4-C5-C6-O6
5	g	3	BMA	C4-C5-C6-O6
5	y	3	BMA	C4-C5-C6-O6
5	z	2	NAG	O5-C5-C6-O6
2	D	1	NAG	C8-C7-N2-C2
2	D	1	NAG	O7-C7-N2-C2
2	K	1	NAG	C8-C7-N2-C2
2	K	1	NAG	O7-C7-N2-C2
2	M	1	NAG	C8-C7-N2-C2
2	M	1	NAG	O7-C7-N2-C2
2	N	1	NAG	C8-C7-N2-C2
2	N	1	NAG	O7-C7-N2-C2
2	U	1	NAG	C8-C7-N2-C2
2	U	1	NAG	O7-C7-N2-C2
2	V	1	NAG	C8-C7-N2-C2
2	V	1	NAG	O7-C7-N2-C2
2	c	1	NAG	C8-C7-N2-C2
2	c	1	NAG	O7-C7-N2-C2
2	e	1	NAG	C8-C7-N2-C2
2	e	1	NAG	O7-C7-N2-C2
2	f	1	NAG	C8-C7-N2-C2
2	f	1	NAG	O7-C7-N2-C2
2	m	1	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	m	1	NAG	O7-C7-N2-C2
2	n	1	NAG	C8-C7-N2-C2
2	n	1	NAG	O7-C7-N2-C2
2	u	1	NAG	C8-C7-N2-C2
2	u	1	NAG	O7-C7-N2-C2
2	w	1	NAG	C8-C7-N2-C2
2	w	1	NAG	O7-C7-N2-C2
2	x	1	NAG	C8-C7-N2-C2
2	x	1	NAG	O7-C7-N2-C2
2	4	1	NAG	C8-C7-N2-C2
2	4	1	NAG	O7-C7-N2-C2
3	G	1	NAG	C8-C7-N2-C2
3	G	1	NAG	O7-C7-N2-C2
3	Y	1	NAG	C8-C7-N2-C2
3	Y	1	NAG	O7-C7-N2-C2
3	q	1	NAG	C8-C7-N2-C2
3	q	1	NAG	O7-C7-N2-C2
5	J	1	NAG	O7-C7-N2-C2
5	L	1	NAG	C8-C7-N2-C2
5	L	1	NAG	O7-C7-N2-C2
5	O	1	NAG	C8-C7-N2-C2
5	O	1	NAG	O7-C7-N2-C2
5	P	1	NAG	C8-C7-N2-C2
5	P	1	NAG	O7-C7-N2-C2
5	b	1	NAG	O7-C7-N2-C2
5	d	1	NAG	C8-C7-N2-C2
5	d	1	NAG	O7-C7-N2-C2
5	g	1	NAG	C8-C7-N2-C2
5	g	1	NAG	O7-C7-N2-C2
5	h	1	NAG	C8-C7-N2-C2
5	h	1	NAG	O7-C7-N2-C2
5	t	1	NAG	O7-C7-N2-C2
5	v	1	NAG	C8-C7-N2-C2
5	v	1	NAG	O7-C7-N2-C2
5	y	1	NAG	C8-C7-N2-C2
5	y	1	NAG	O7-C7-N2-C2
5	z	1	NAG	C8-C7-N2-C2
5	z	1	NAG	O7-C7-N2-C2
6	R	1	NAG	C8-C7-N2-C2
6	R	1	NAG	O7-C7-N2-C2
6	j	1	NAG	C8-C7-N2-C2
6	j	1	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
6	1	1	NAG	C8-C7-N2-C2
6	1	1	NAG	O7-C7-N2-C2
7	S	1	NAG	C8-C7-N2-C2
7	S	1	NAG	O7-C7-N2-C2
7	k	1	NAG	C8-C7-N2-C2
7	k	1	NAG	O7-C7-N2-C2
7	2	1	NAG	C8-C7-N2-C2
7	2	1	NAG	O7-C7-N2-C2
5	P	2	NAG	O5-C5-C6-O6
5	h	2	NAG	O5-C5-C6-O6
8	T	4	MAN	O5-C5-C6-O6
8	l	4	MAN	O5-C5-C6-O6
8	3	4	MAN	O5-C5-C6-O6
5	J	1	NAG	C4-C5-C6-O6
5	b	1	NAG	C4-C5-C6-O6
5	t	1	NAG	C4-C5-C6-O6
2	M	2	NAG	O5-C5-C6-O6
2	e	2	NAG	O5-C5-C6-O6
2	w	2	NAG	O5-C5-C6-O6
4	I	3	BMA	O5-C5-C6-O6
4	a	3	BMA	O5-C5-C6-O6
4	s	3	BMA	O5-C5-C6-O6
5	J	1	NAG	O5-C5-C6-O6
5	b	1	NAG	O5-C5-C6-O6
5	t	1	NAG	O5-C5-C6-O6
2	x	1	NAG	C4-C5-C6-O6
2	N	1	NAG	O5-C5-C6-O6
2	x	1	NAG	O5-C5-C6-O6
2	f	1	NAG	C4-C5-C6-O6
2	f	1	NAG	O5-C5-C6-O6
2	N	1	NAG	C4-C5-C6-O6
8	T	7	MAN	C4-C5-C6-O6
8	l	7	MAN	C4-C5-C6-O6
8	3	7	MAN	C4-C5-C6-O6
2	N	2	NAG	C4-C5-C6-O6
2	f	2	NAG	C4-C5-C6-O6
2	x	2	NAG	C4-C5-C6-O6
3	G	3	BMA	O5-C5-C6-O6
3	Y	3	BMA	O5-C5-C6-O6
3	q	3	BMA	O5-C5-C6-O6
4	I	5	MAN	O5-C5-C6-O6
4	a	5	MAN	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	s	5	MAN	O5-C5-C6-O6
5	L	4	MAN	O5-C5-C6-O6
5	d	4	MAN	O5-C5-C6-O6
5	v	4	MAN	O5-C5-C6-O6
5	O	2	NAG	O5-C5-C6-O6
5	g	2	NAG	O5-C5-C6-O6
5	y	2	NAG	O5-C5-C6-O6
5	L	2	NAG	O5-C5-C6-O6
5	d	2	NAG	O5-C5-C6-O6
5	v	2	NAG	O5-C5-C6-O6
8	T	7	MAN	O5-C5-C6-O6
8	l	7	MAN	O5-C5-C6-O6
8	3	7	MAN	O5-C5-C6-O6
2	m	1	NAG	C4-C5-C6-O6
2	4	1	NAG	C4-C5-C6-O6
2	U	1	NAG	C4-C5-C6-O6
5	P	2	NAG	C4-C5-C6-O6
5	h	2	NAG	C4-C5-C6-O6
5	z	2	NAG	C4-C5-C6-O6
2	0	2	NAG	C4-C5-C6-O6
2	Q	2	NAG	C4-C5-C6-O6
2	i	2	NAG	C4-C5-C6-O6
5	d	1	NAG	C4-C5-C6-O6
5	L	1	NAG	C4-C5-C6-O6
5	v	1	NAG	C4-C5-C6-O6
2	4	1	NAG	O5-C5-C6-O6
2	m	1	NAG	O5-C5-C6-O6
2	U	1	NAG	O5-C5-C6-O6
2	D	2	NAG	C3-C2-N2-C7
2	F	1	NAG	C3-C2-N2-C7
2	V	2	NAG	C3-C2-N2-C7
2	X	1	NAG	C3-C2-N2-C7
2	n	2	NAG	C3-C2-N2-C7
2	p	1	NAG	C3-C2-N2-C7
8	T	1	NAG	C3-C2-N2-C7
8	T	2	NAG	C3-C2-N2-C7
8	l	1	NAG	C3-C2-N2-C7
8	l	2	NAG	C3-C2-N2-C7
8	3	1	NAG	C3-C2-N2-C7
8	3	2	NAG	C3-C2-N2-C7
5	J	3	BMA	O5-C5-C6-O6
5	t	3	BMA	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	b	3	BMA	O5-C5-C6-O6
7	S	1	NAG	C4-C5-C6-O6
7	k	1	NAG	C4-C5-C6-O6
7	2	1	NAG	C4-C5-C6-O6
2	F	1	NAG	C1-C2-N2-C7
2	X	1	NAG	C1-C2-N2-C7
2	p	1	NAG	C1-C2-N2-C7
8	T	2	NAG	C1-C2-N2-C7
8	l	2	NAG	C1-C2-N2-C7
8	3	2	NAG	C1-C2-N2-C7
3	E	1	NAG	C3-C2-N2-C7
3	W	1	NAG	C3-C2-N2-C7
3	o	1	NAG	C3-C2-N2-C7
8	3	2	NAG	C4-C5-C6-O6
8	l	2	NAG	C4-C5-C6-O6
8	T	2	NAG	C4-C5-C6-O6
2	i	2	NAG	O5-C5-C6-O6
2	Q	2	NAG	O5-C5-C6-O6
2	0	2	NAG	O5-C5-C6-O6
8	l	3	BMA	C4-C5-C6-O6
8	T	3	BMA	C4-C5-C6-O6
8	3	3	BMA	C4-C5-C6-O6
5	L	1	NAG	O5-C5-C6-O6
5	d	1	NAG	O5-C5-C6-O6
5	v	1	NAG	O5-C5-C6-O6

There are no ring outliers.

70 monomers are involved in 113 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	V	2	NAG	1	0
5	t	4	MAN	2	0
5	b	2	NAG	1	0
3	Z	3	BMA	1	0
5	g	2	NAG	2	0
5	J	2	NAG	1	0
5	y	2	NAG	1	0
2	w	1	NAG	2	0
3	H	1	NAG	5	0
5	J	3	BMA	2	0
5	y	3	BMA	2	0
2	F	1	NAG	4	0

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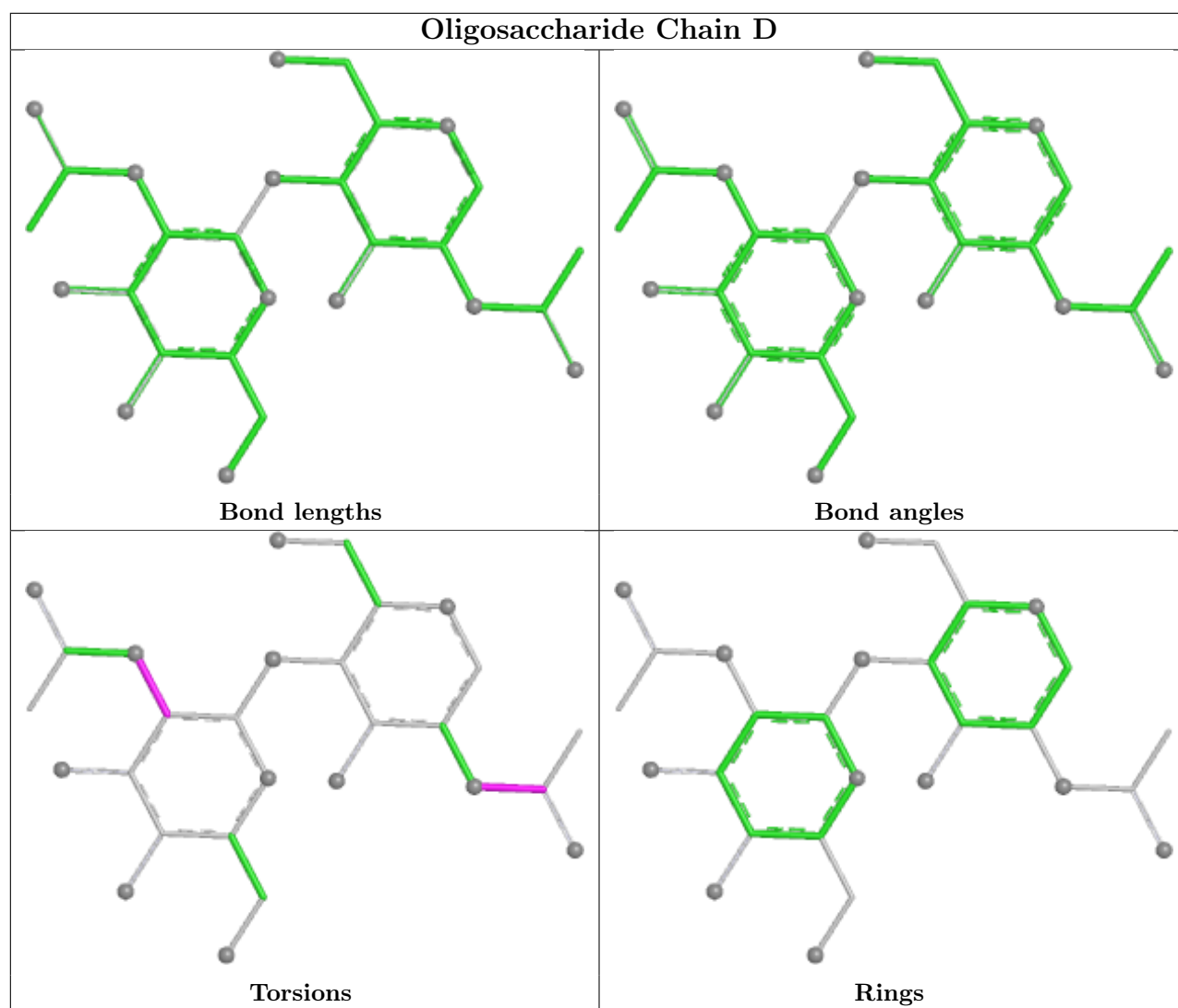
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	e	1	NAG	1	0
8	T	1	NAG	3	0
2	n	1	NAG	1	0
8	l	1	NAG	3	0
4	I	1	NAG	2	0
3	G	2	NAG	5	0
5	t	2	NAG	1	0
2	D	1	NAG	2	0
6	j	4	FUC	3	0
5	g	1	NAG	1	0
5	b	1	NAG	1	0
2	D	2	NAG	1	0
6	l	4	FUC	4	0
8	T	6	MAN	1	0
4	s	1	NAG	2	0
3	r	1	NAG	4	0
5	g	3	BMA	2	0
5	O	1	NAG	1	0
2	n	2	NAG	1	0
2	X	1	NAG	3	0
3	r	3	BMA	1	0
2	M	1	NAG	2	0
6	R	4	FUC	4	0
5	J	4	MAN	2	0
3	Z	1	NAG	5	0
8	3	1	NAG	3	0
4	a	1	NAG	2	0
8	3	7	MAN	1	0
3	Z	2	NAG	2	0
3	r	2	NAG	2	0
6	l	2	NAG	4	0
6	j	2	NAG	3	0
5	y	1	NAG	1	0
5	J	1	NAG	1	0
8	3	6	MAN	1	0
4	a	5	MAN	1	0
5	b	3	BMA	2	0
5	t	3	BMA	2	0
2	e	2	NAG	1	0
3	Y	2	NAG	5	0
8	l	7	MAN	1	0
2	p	1	NAG	3	0

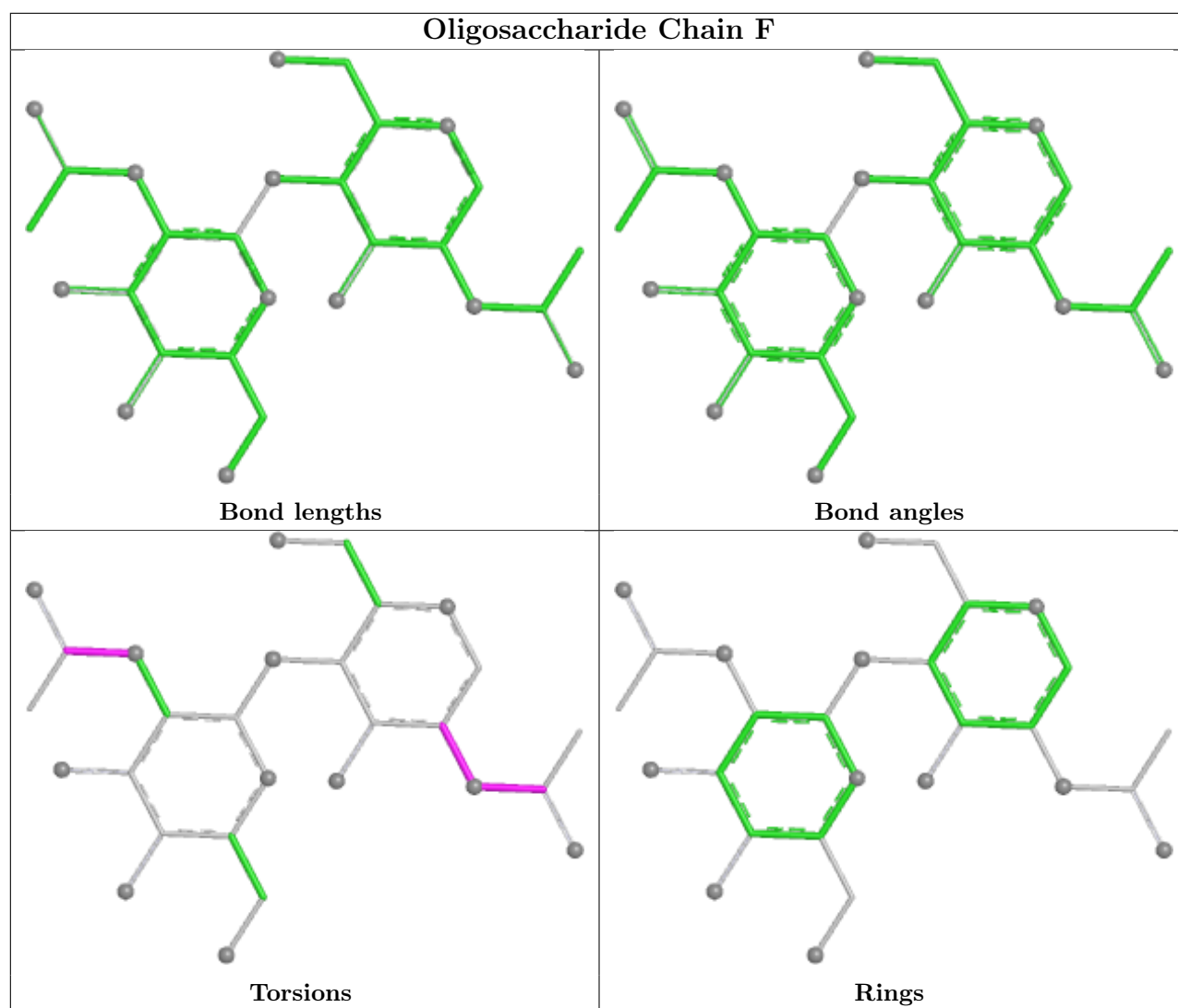
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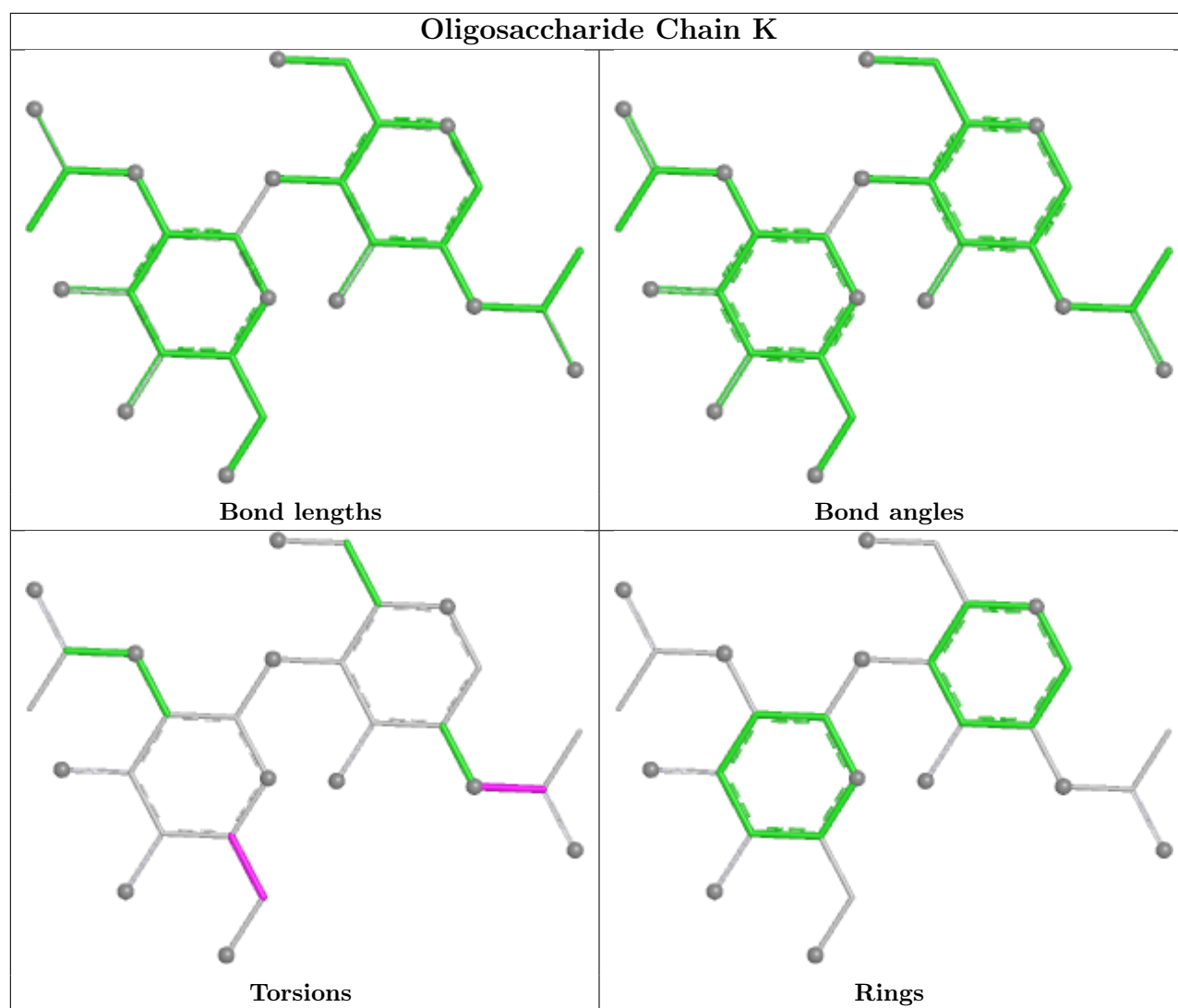
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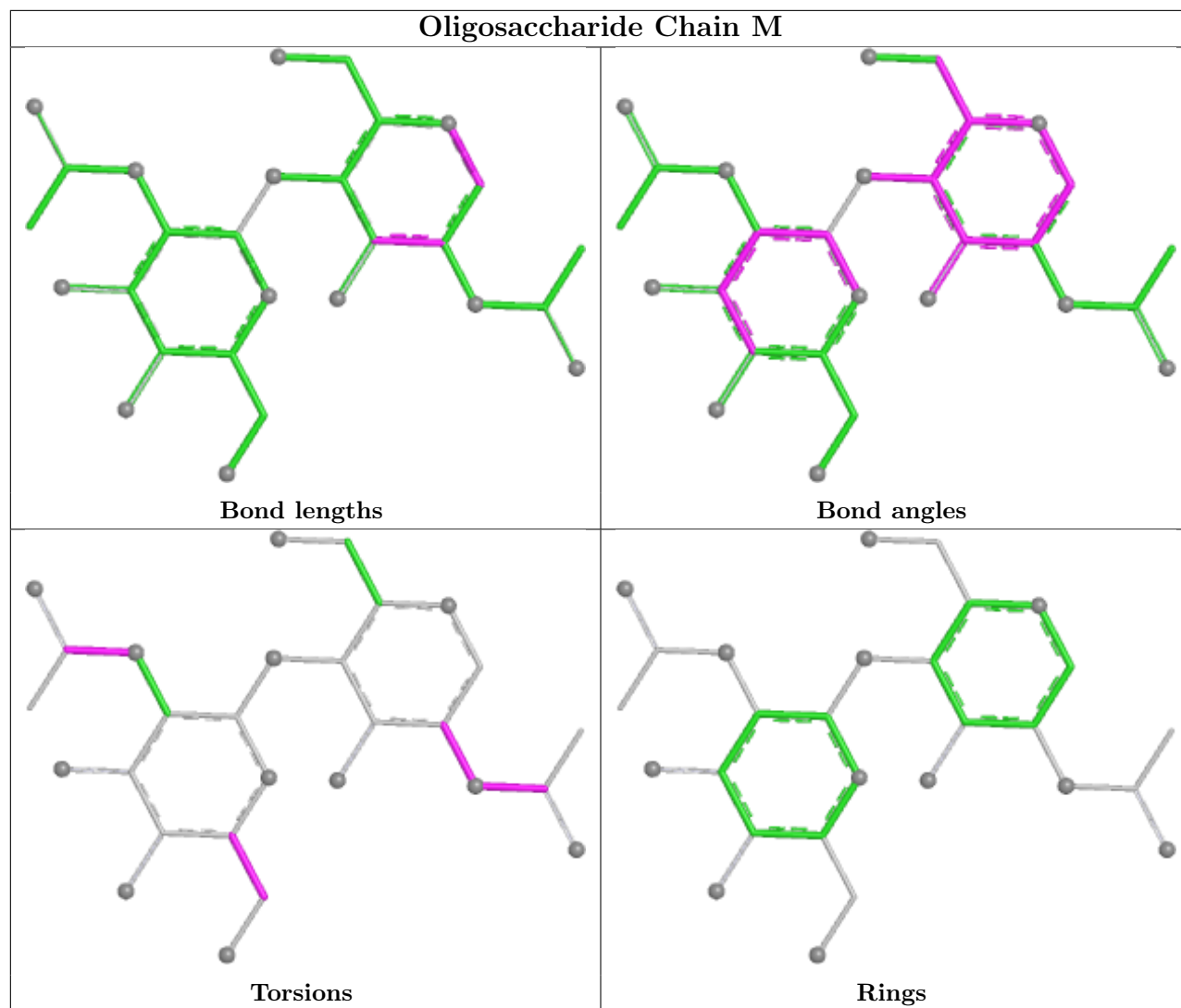
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	s	5	MAN	1	0
3	G	1	NAG	3	0
5	b	4	MAN	2	0
6	R	2	NAG	4	0
3	H	3	BMA	1	0
4	I	5	MAN	1	0
3	q	1	NAG	3	0
8	l	6	MAN	1	0
3	H	2	NAG	2	0
5	O	3	BMA	1	0
2	V	1	NAG	1	0
2	M	2	NAG	1	0
3	Y	1	NAG	3	0
8	T	7	MAN	1	0
5	t	1	NAG	2	0
3	q	2	NAG	5	0

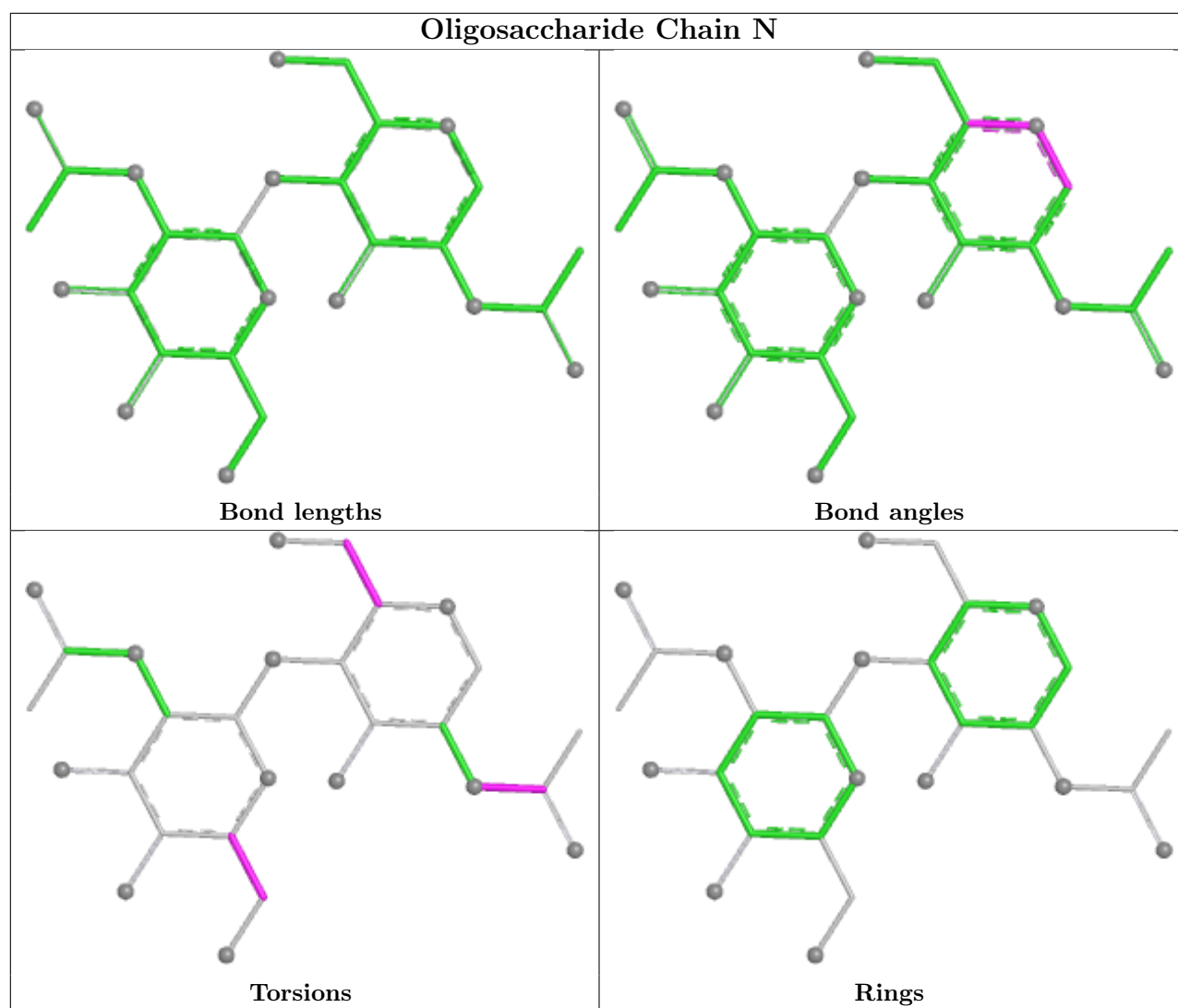
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

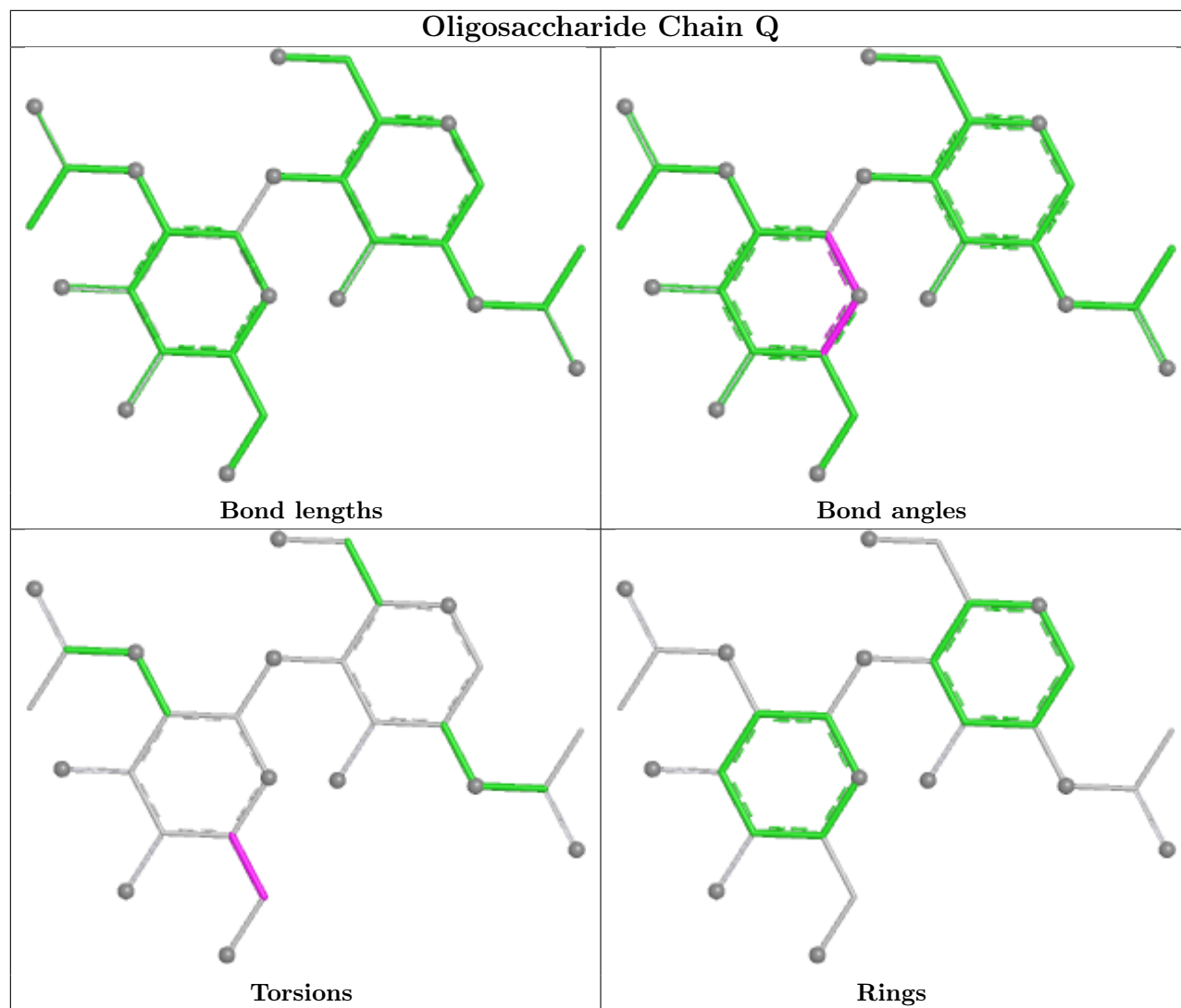


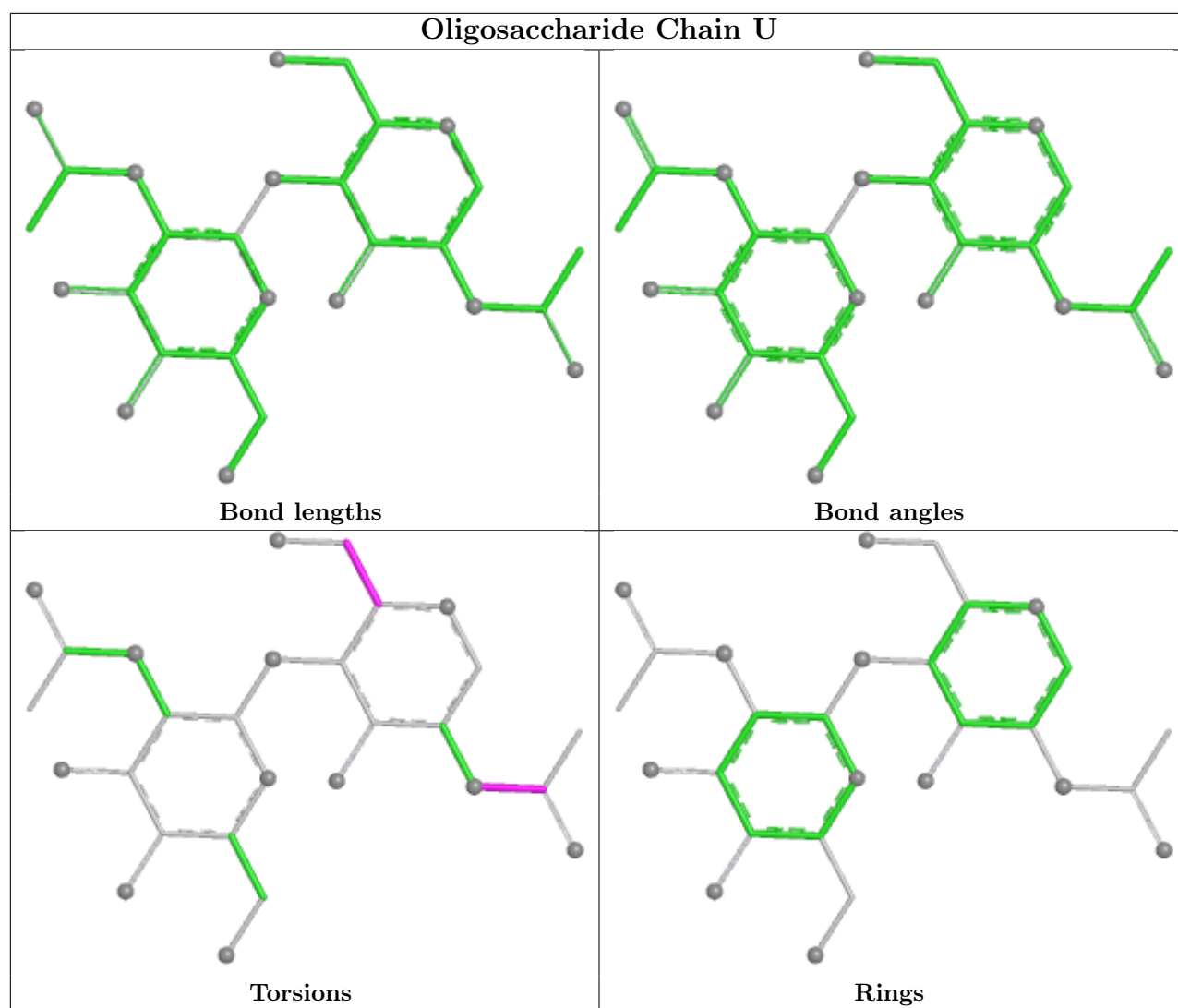


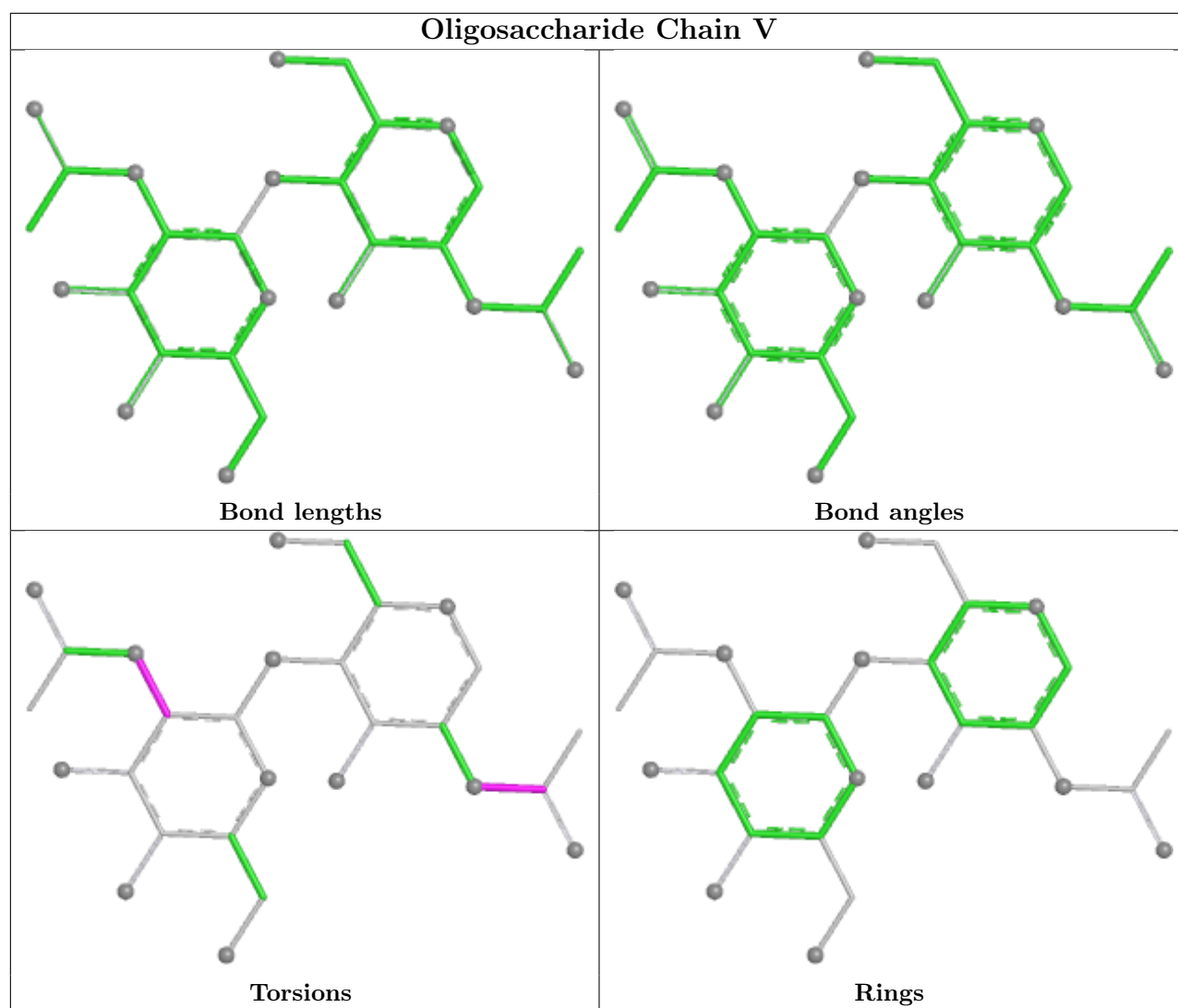


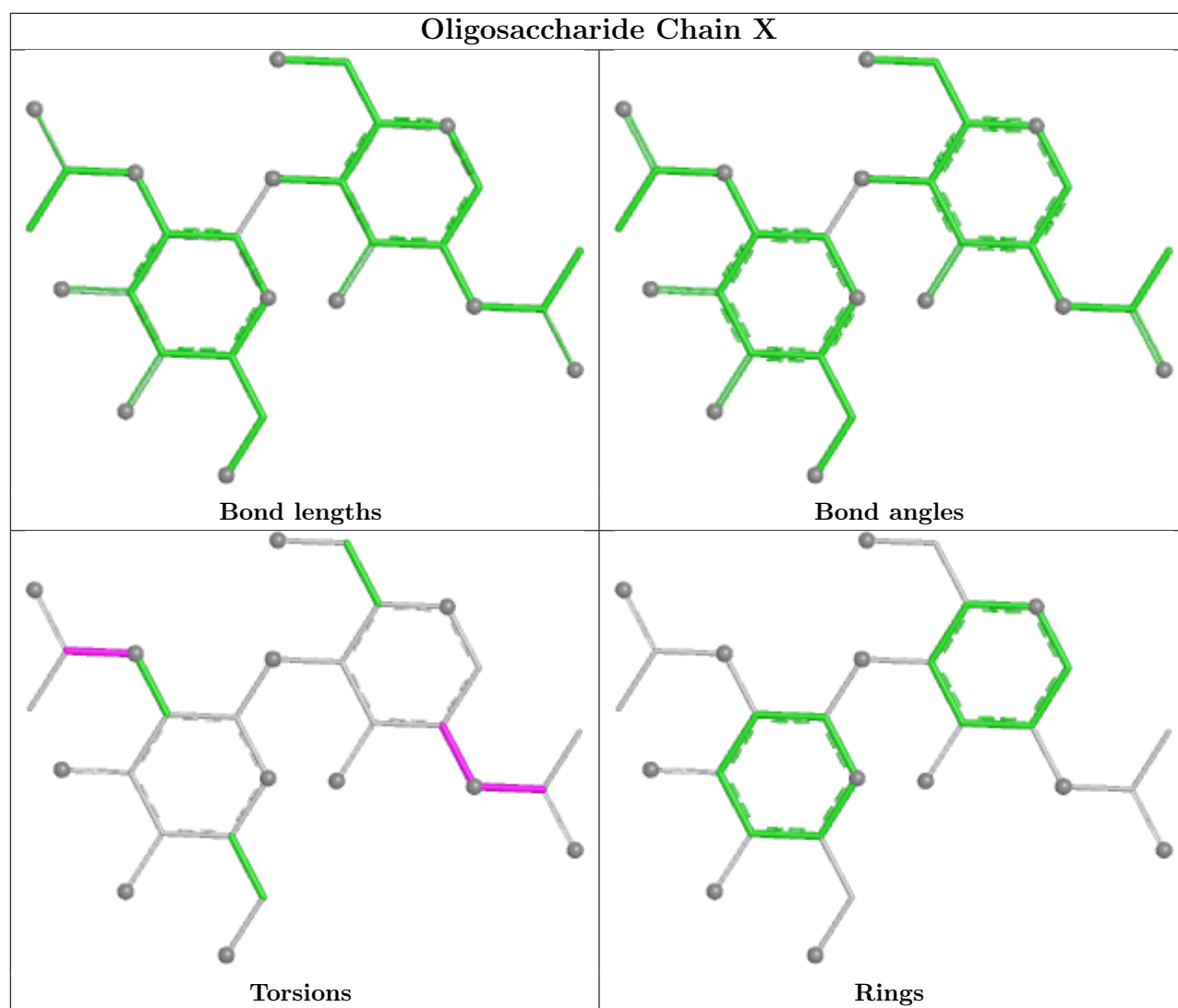


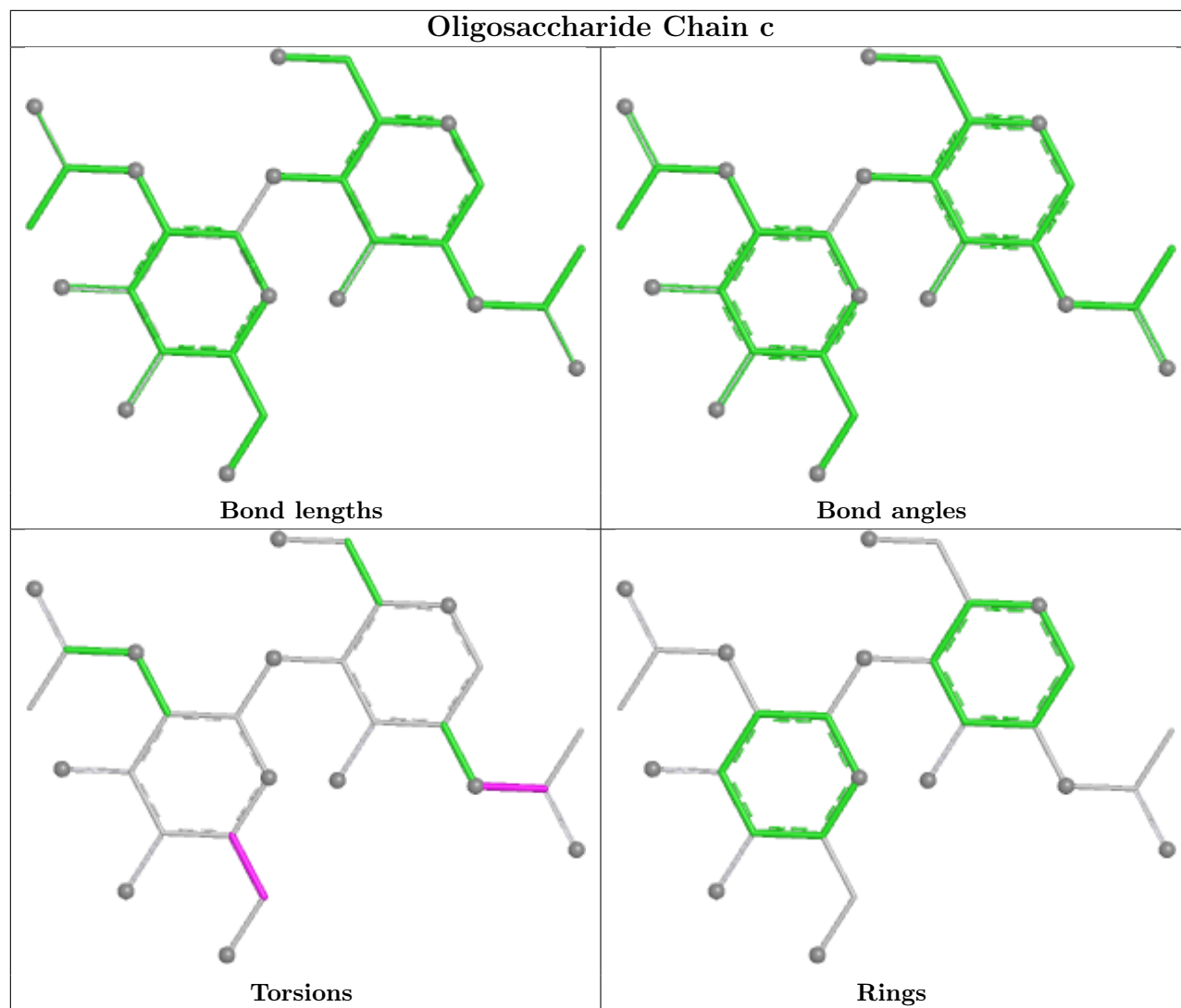


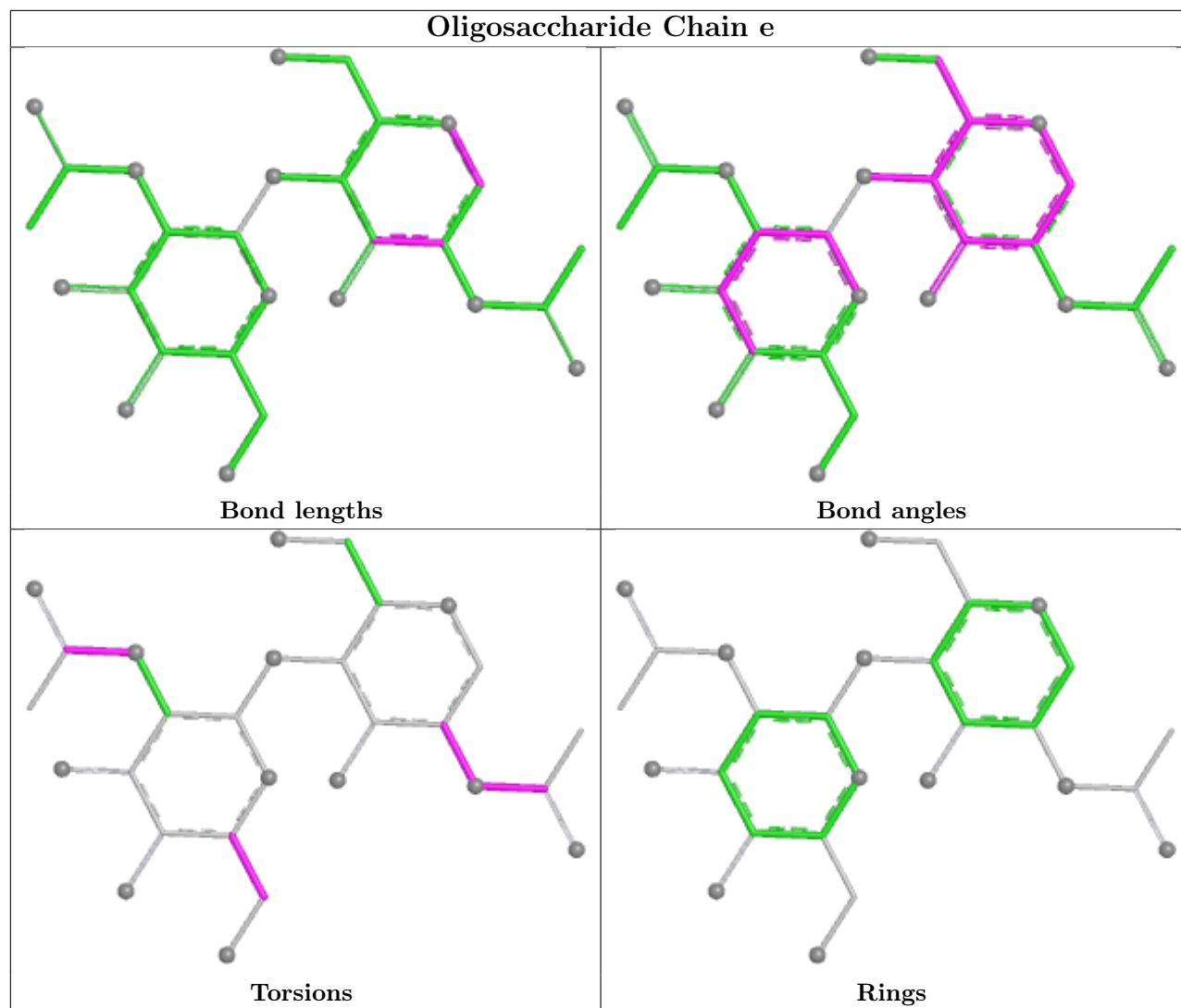


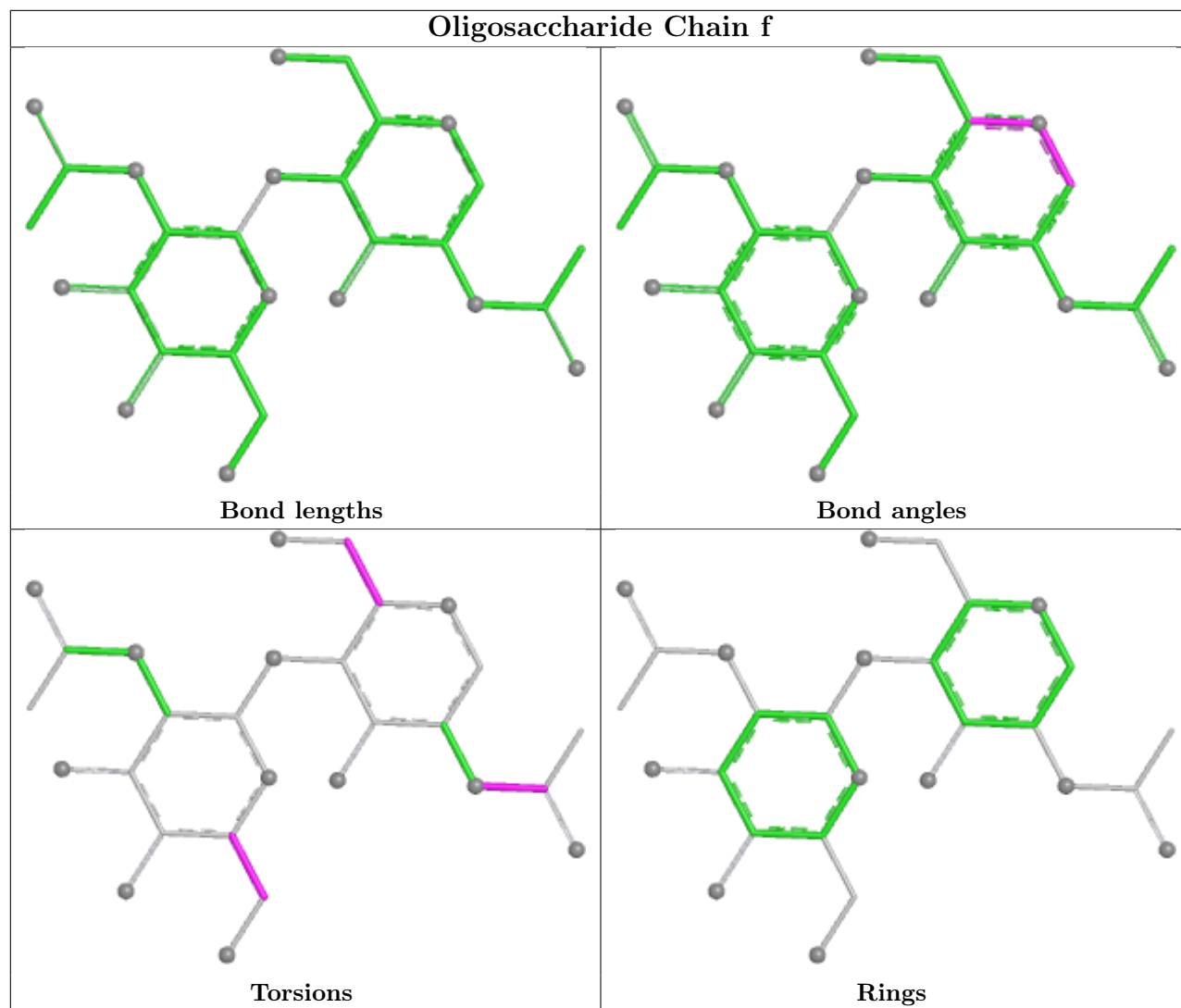


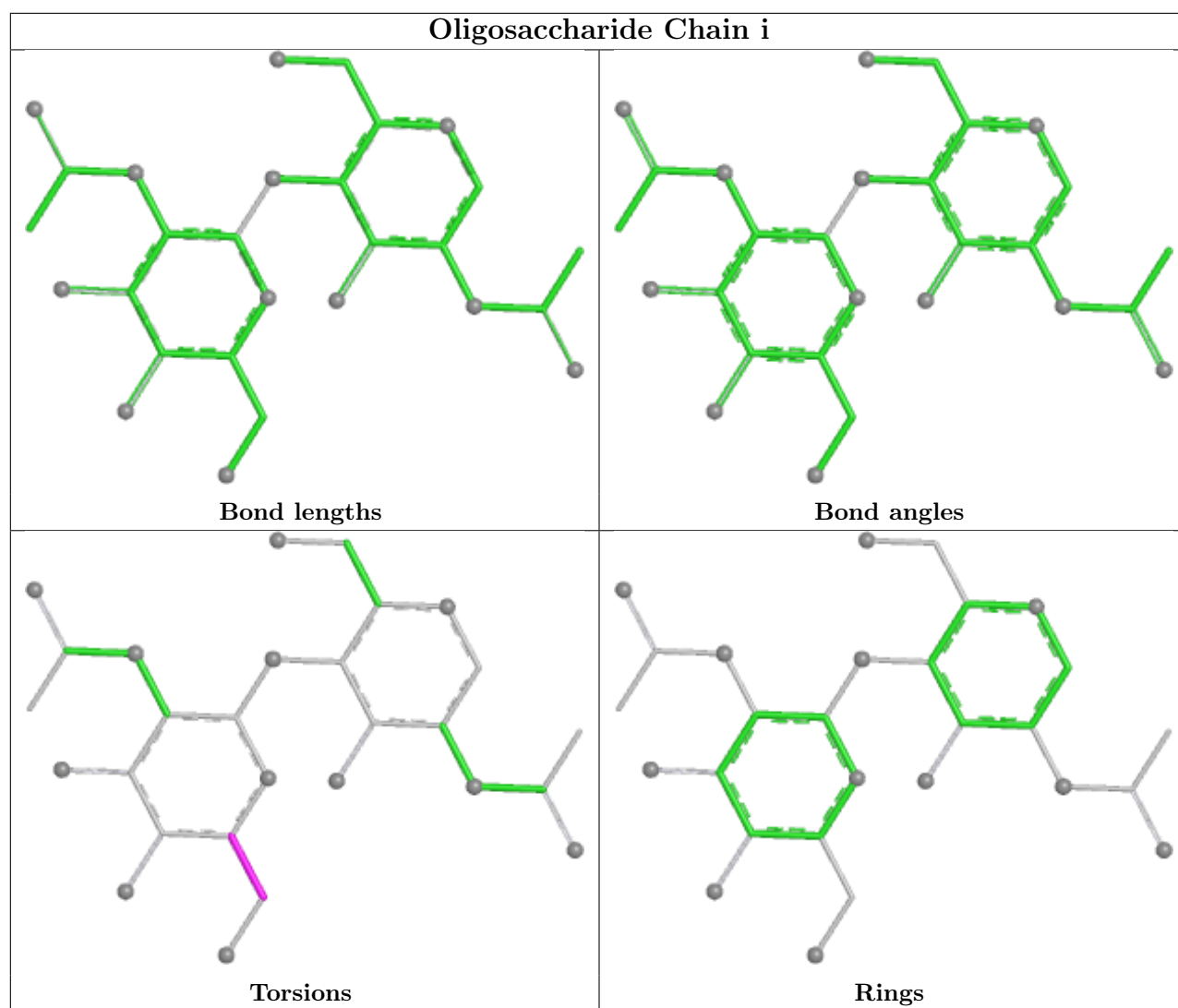


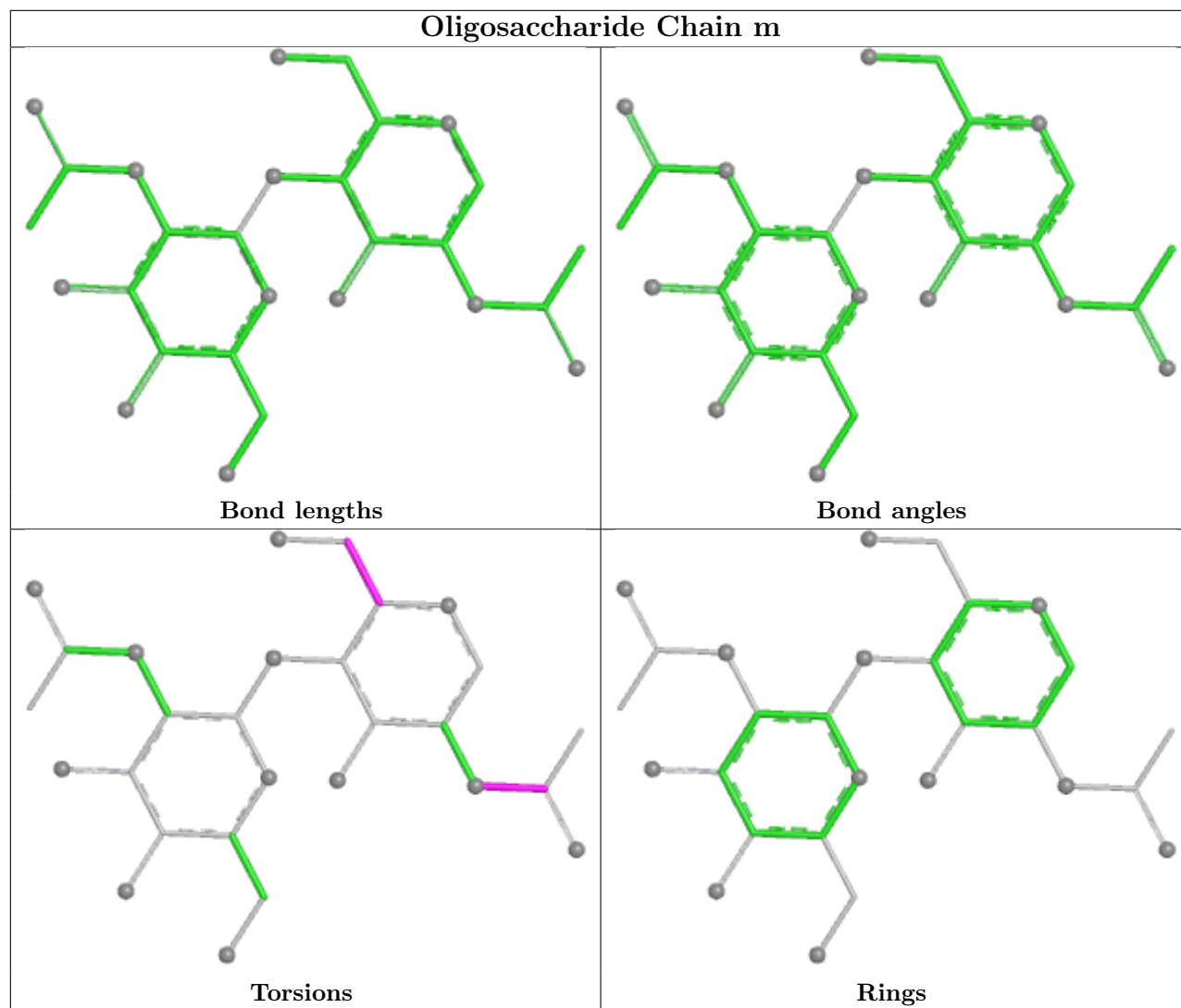


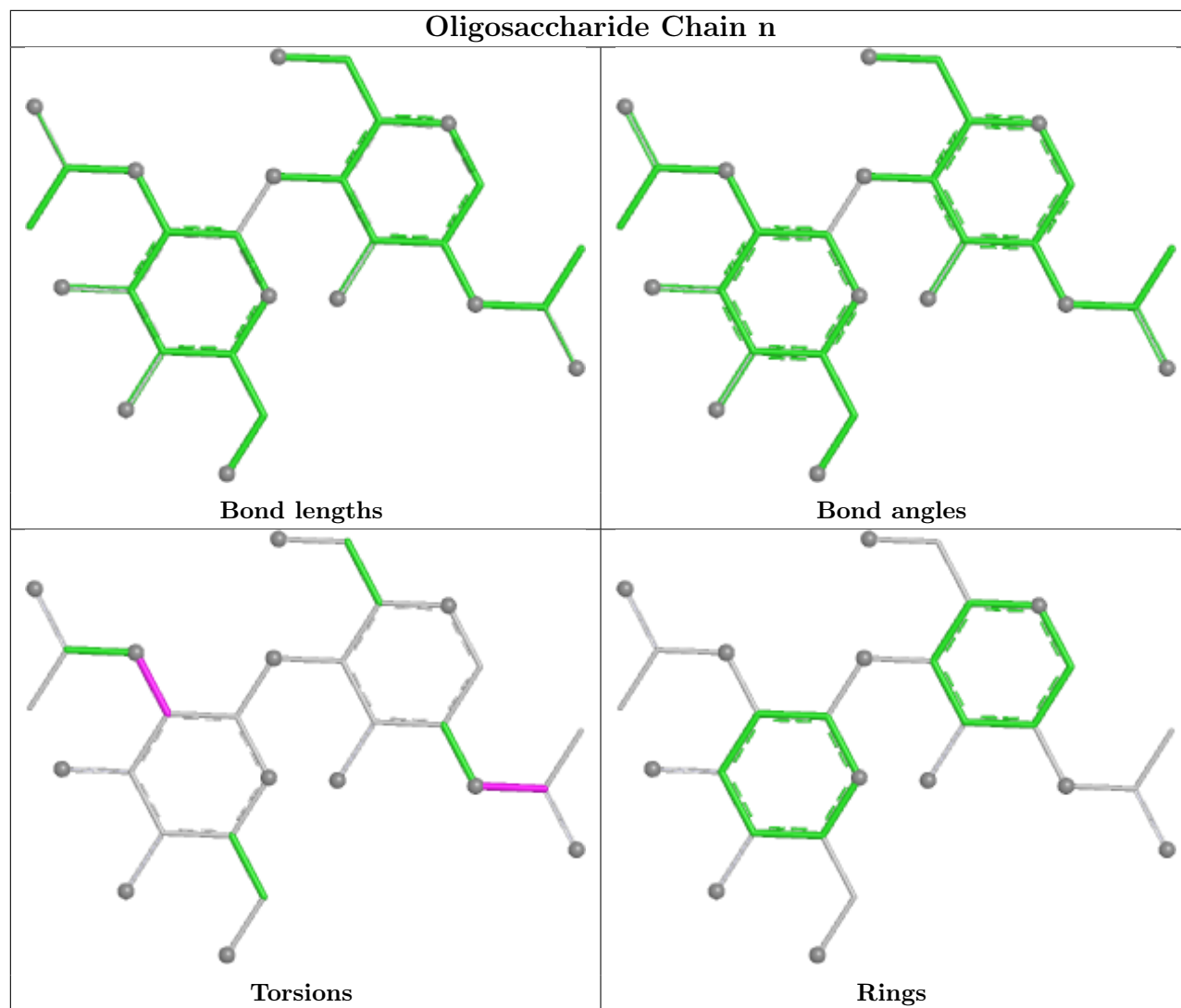


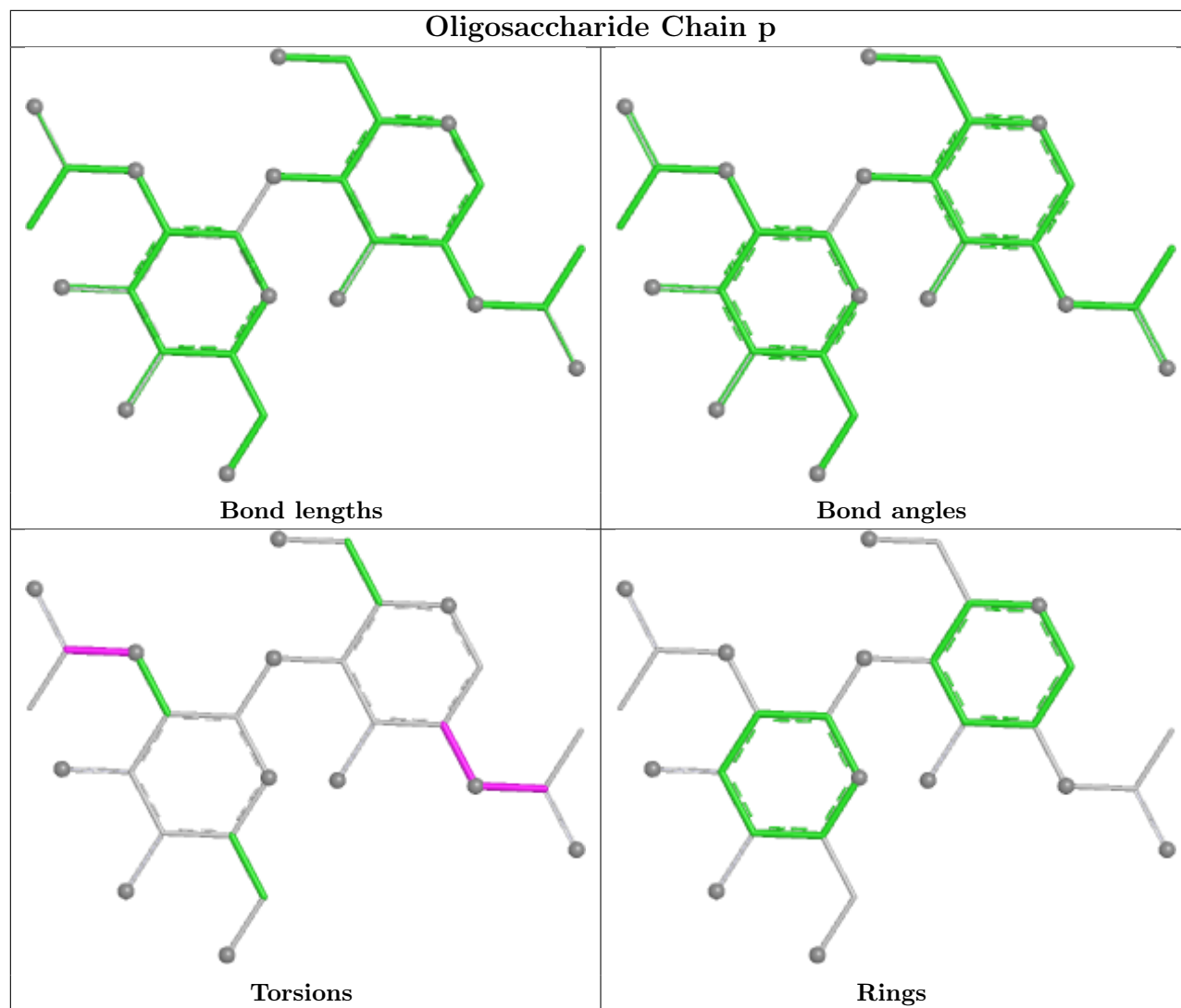


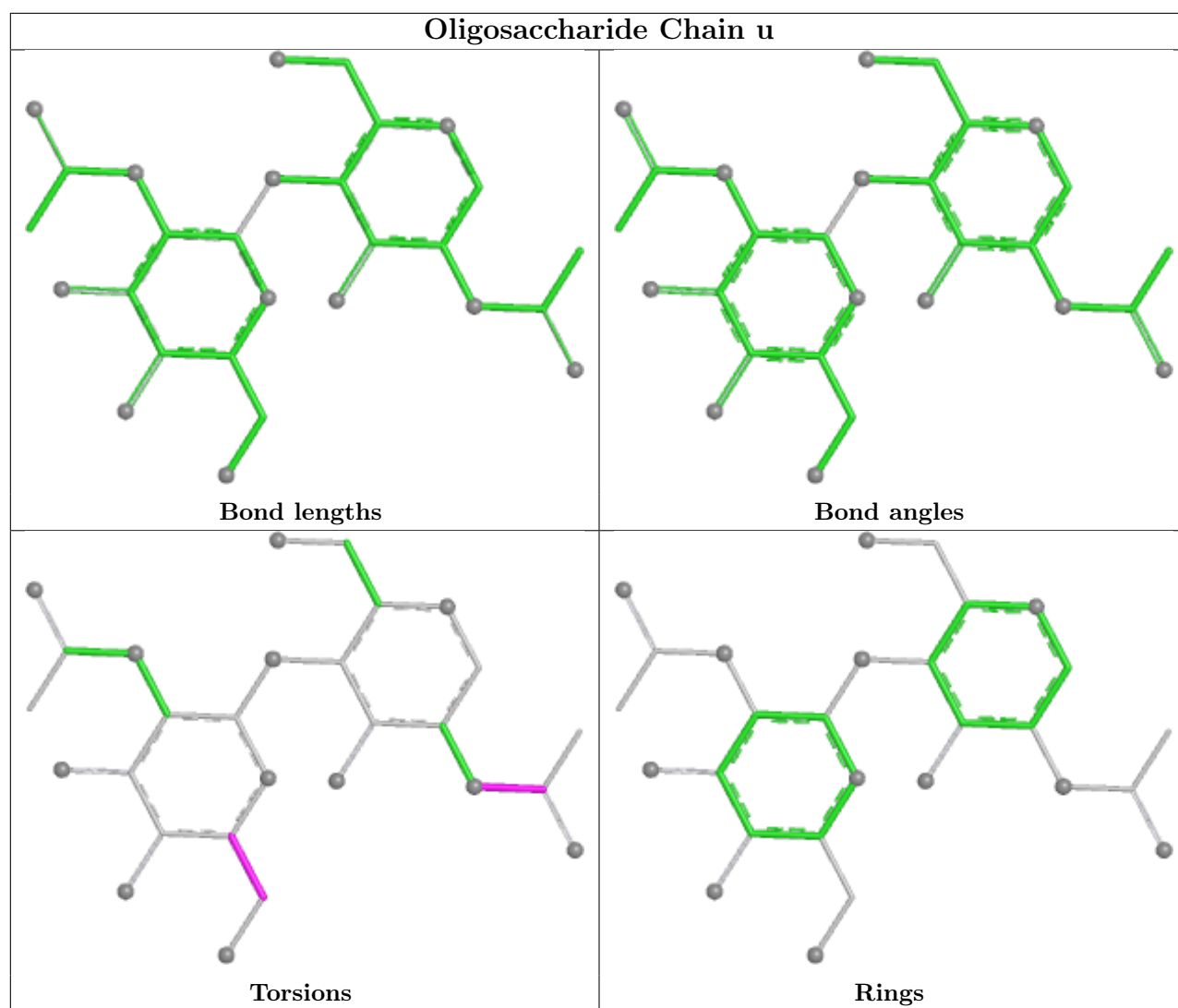


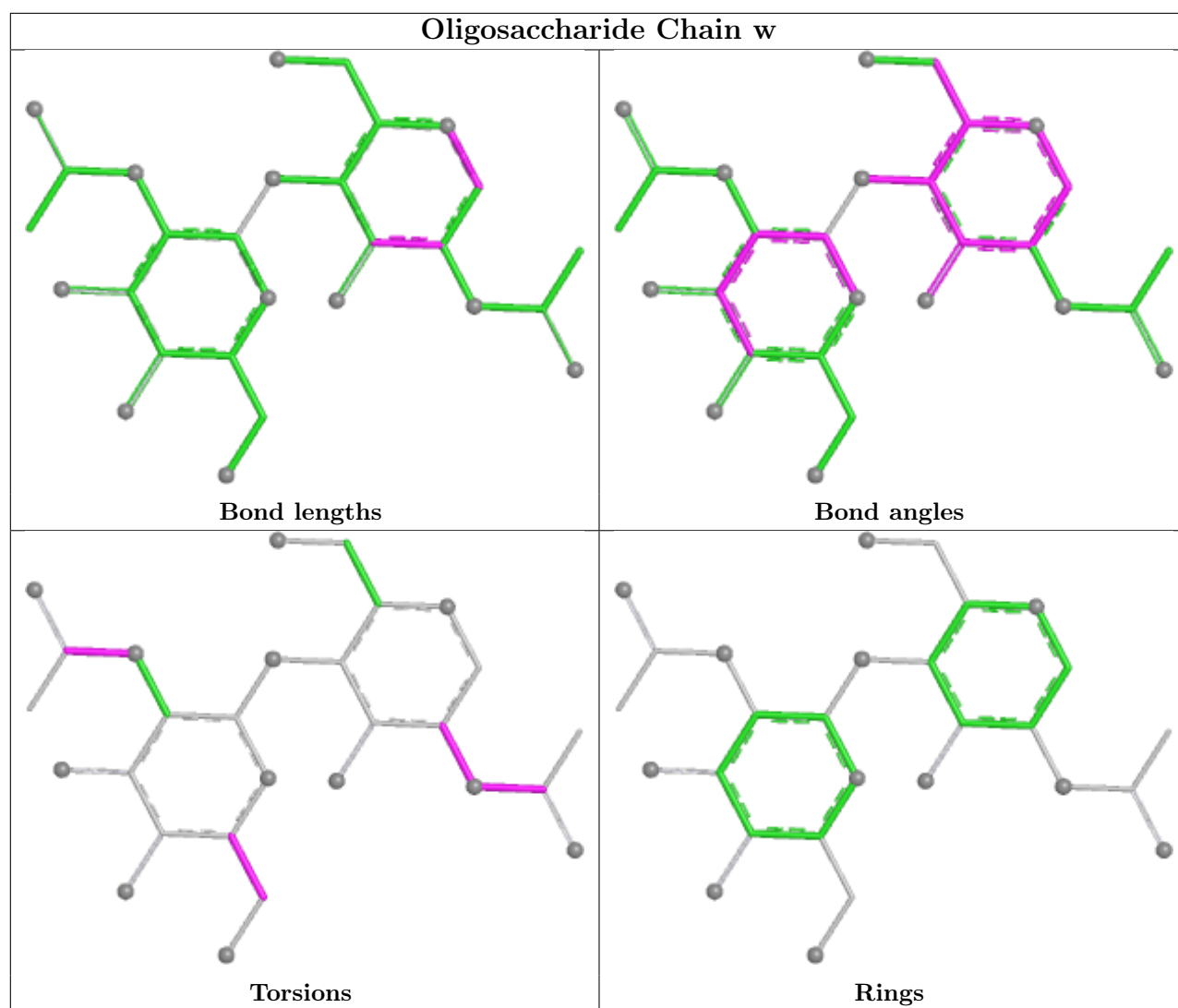


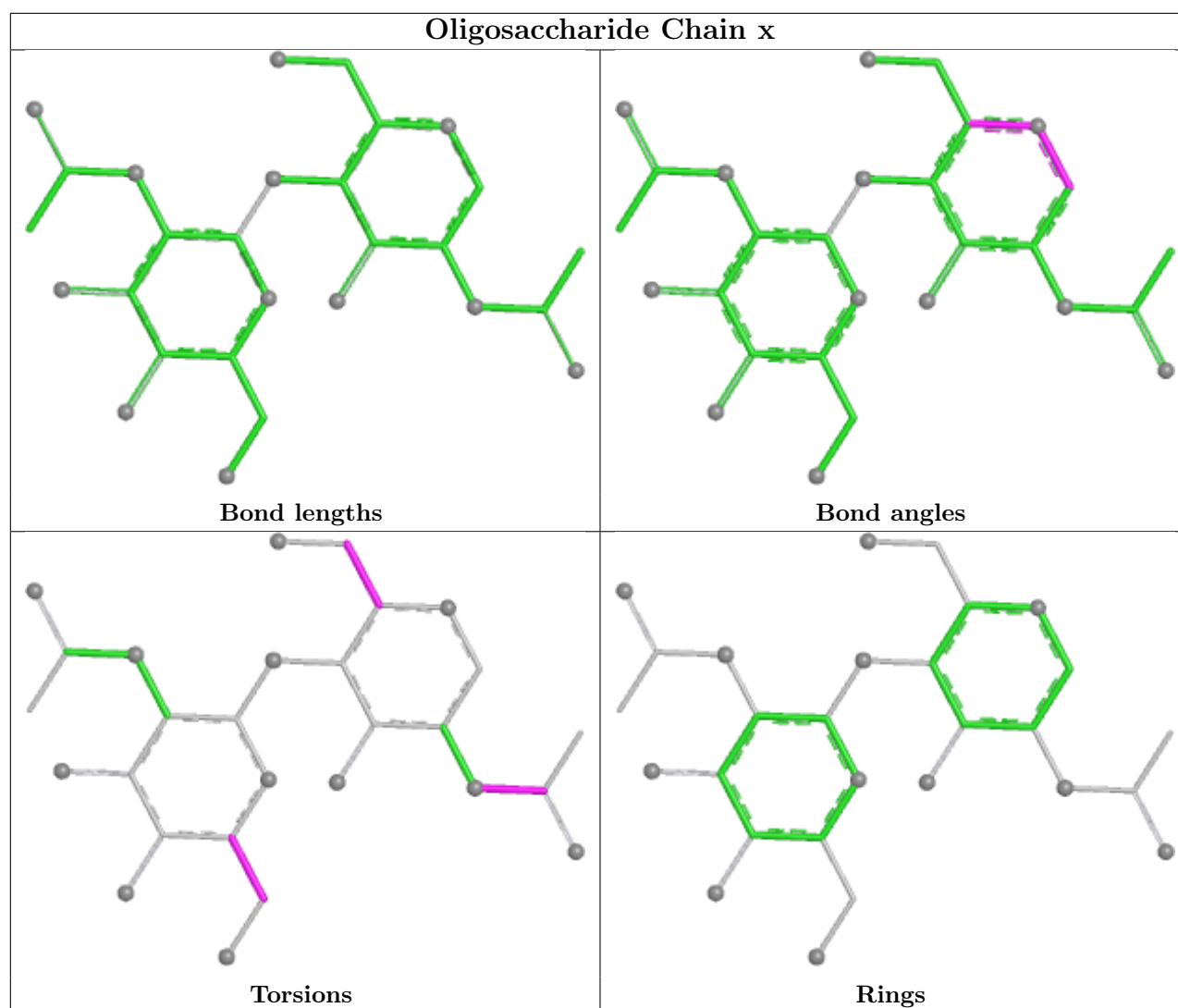


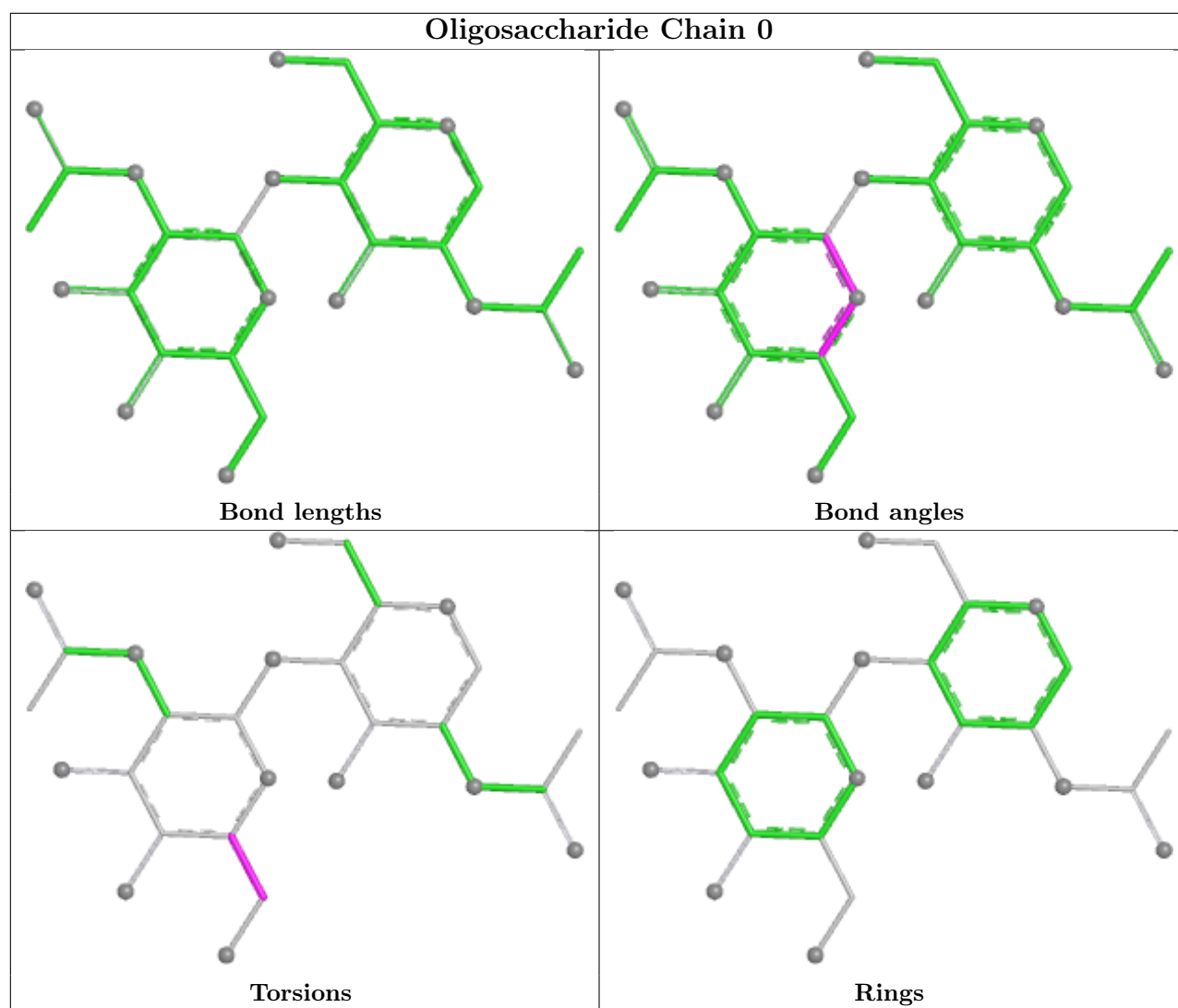


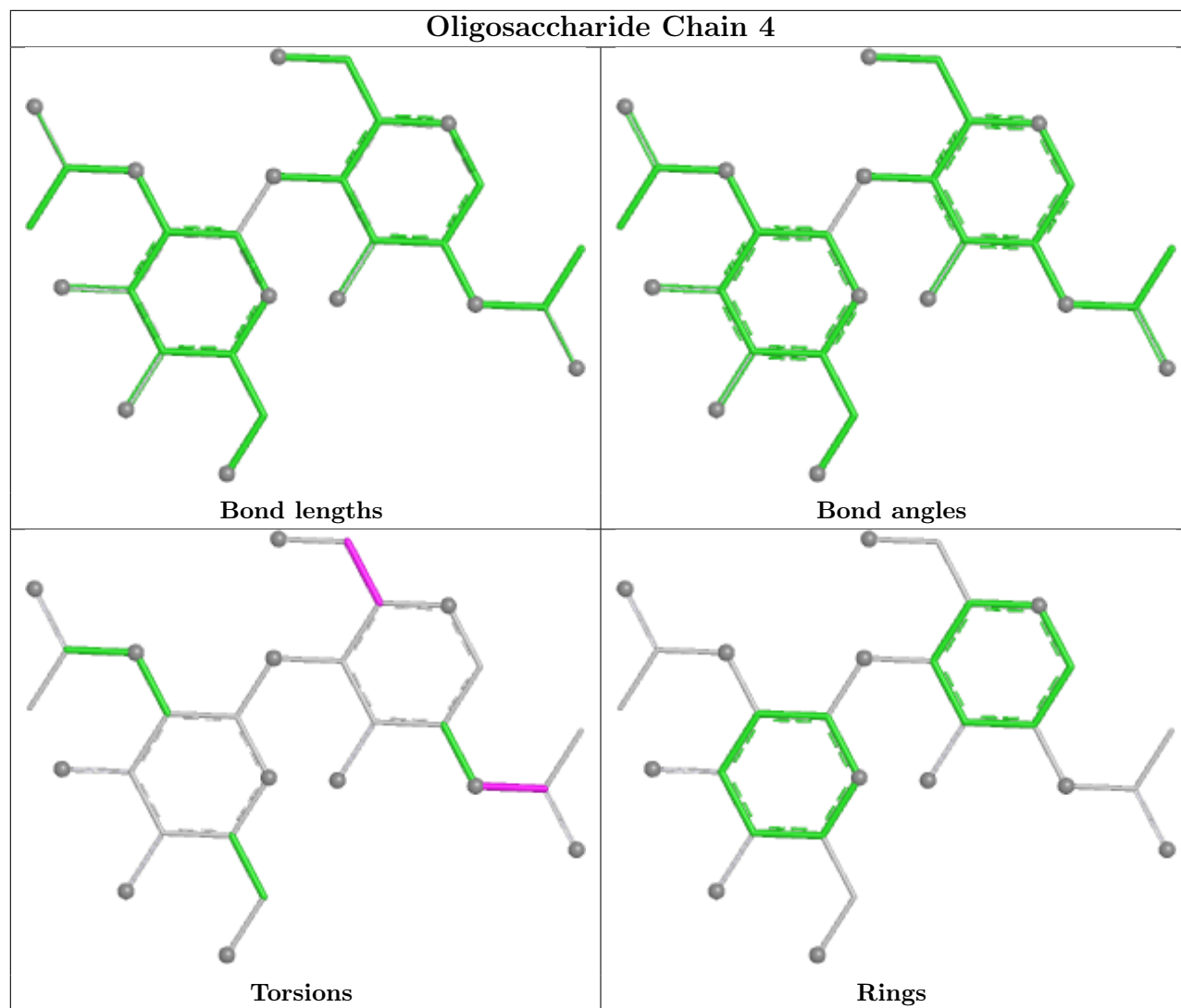


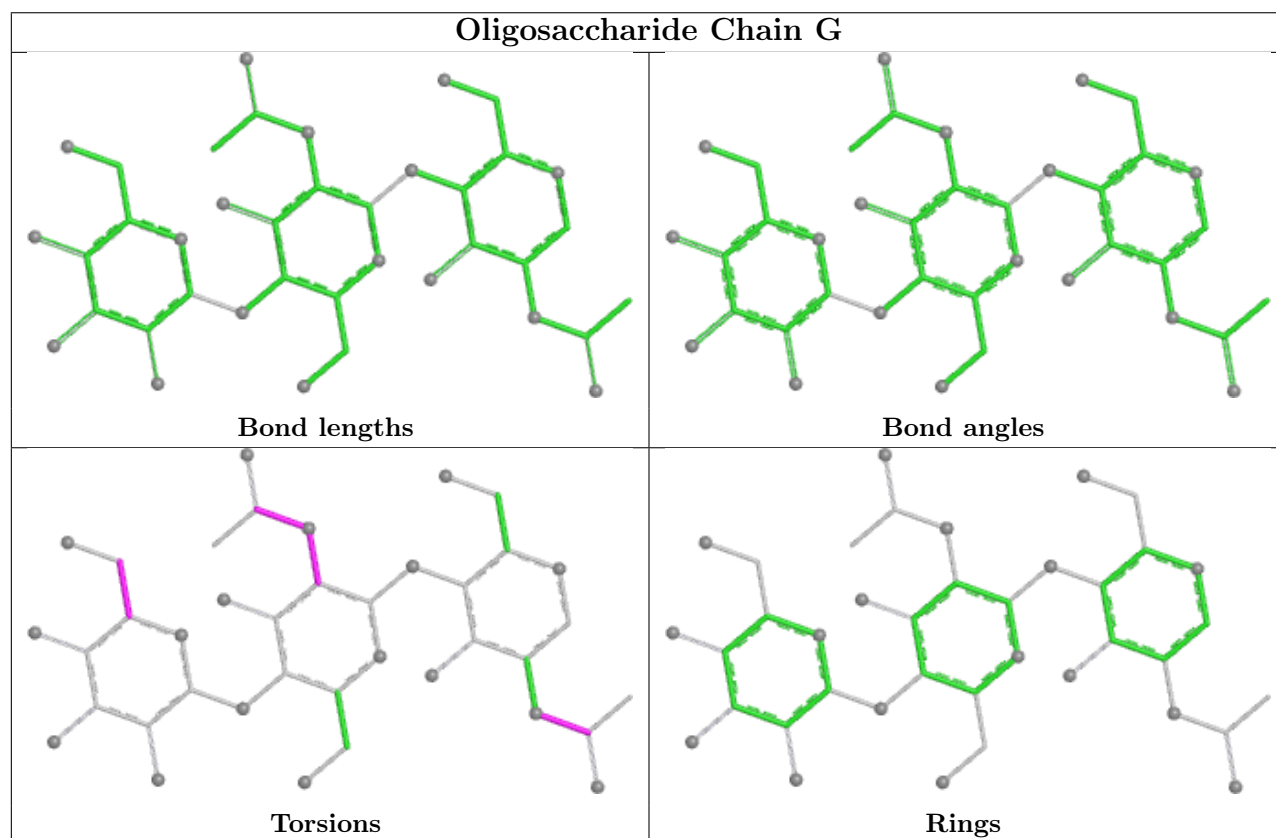
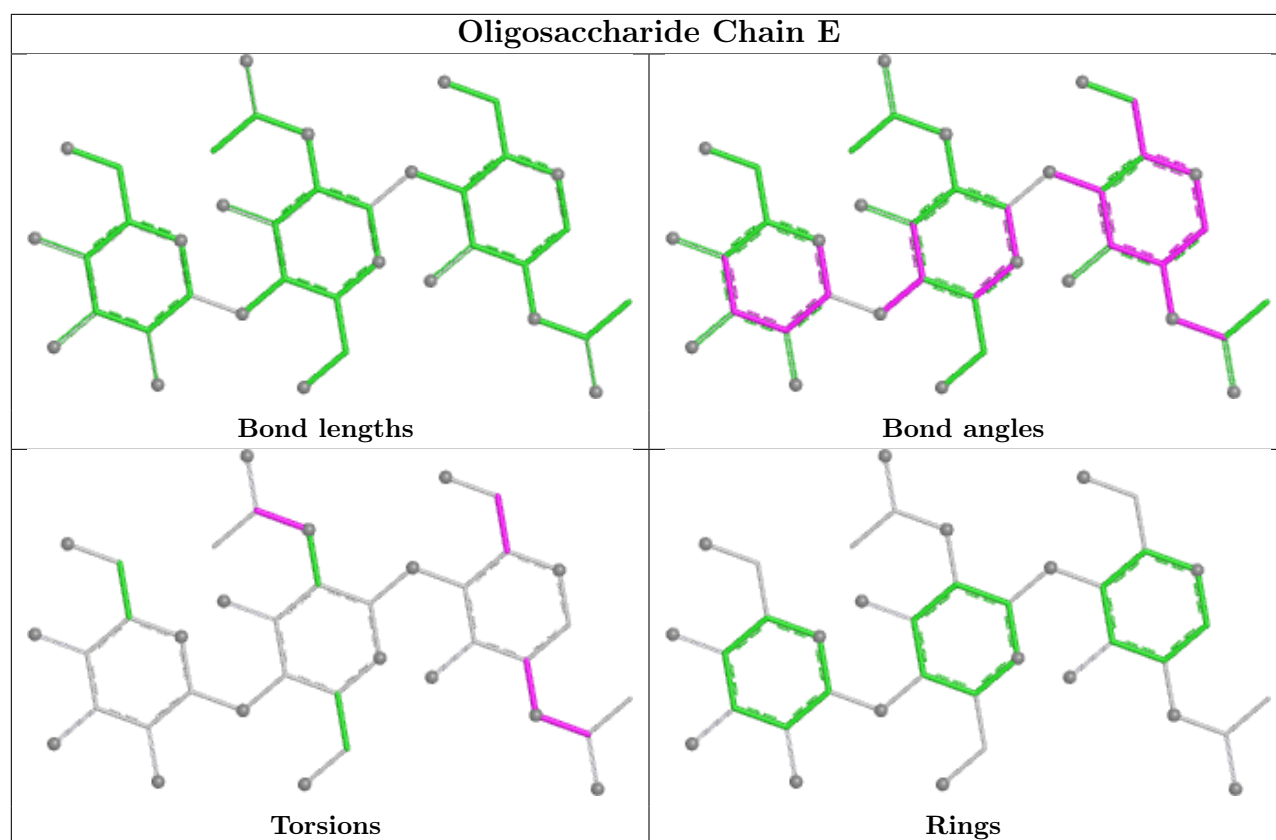


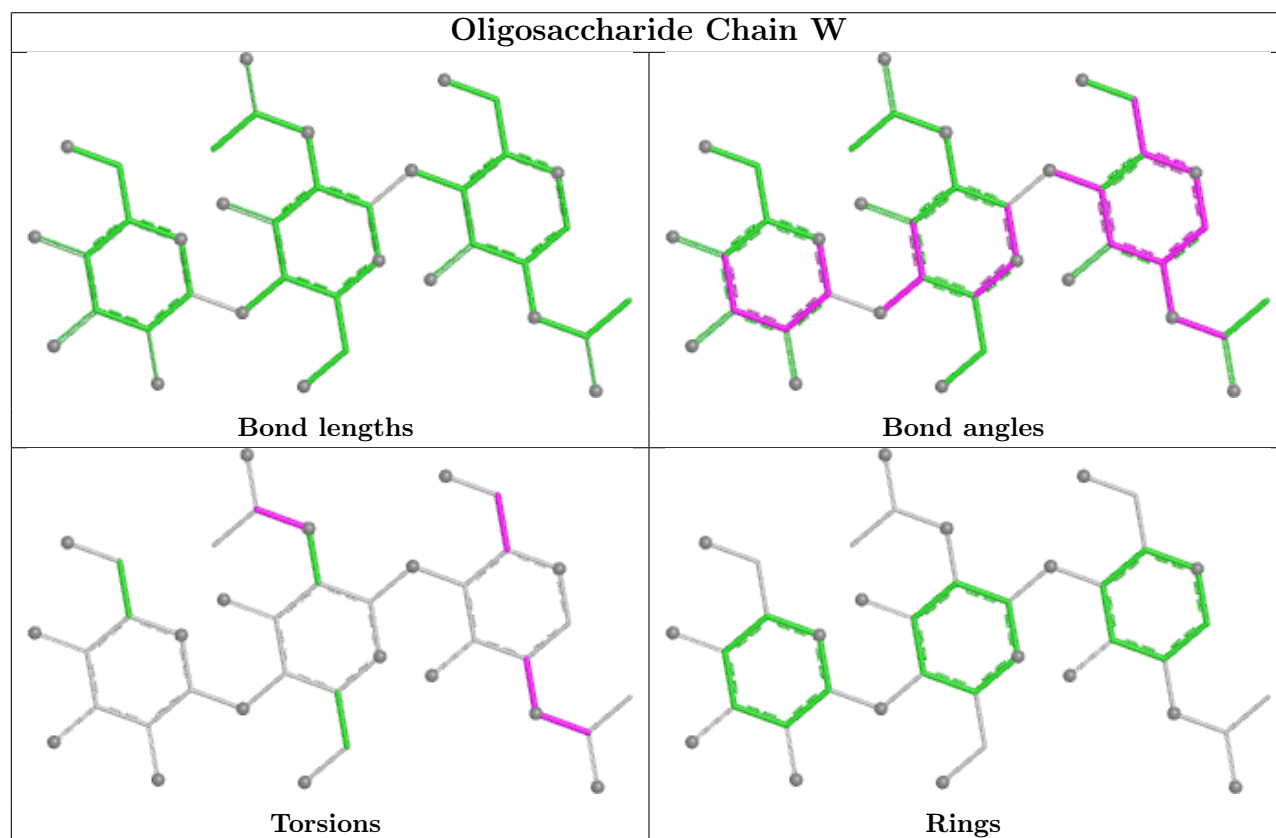
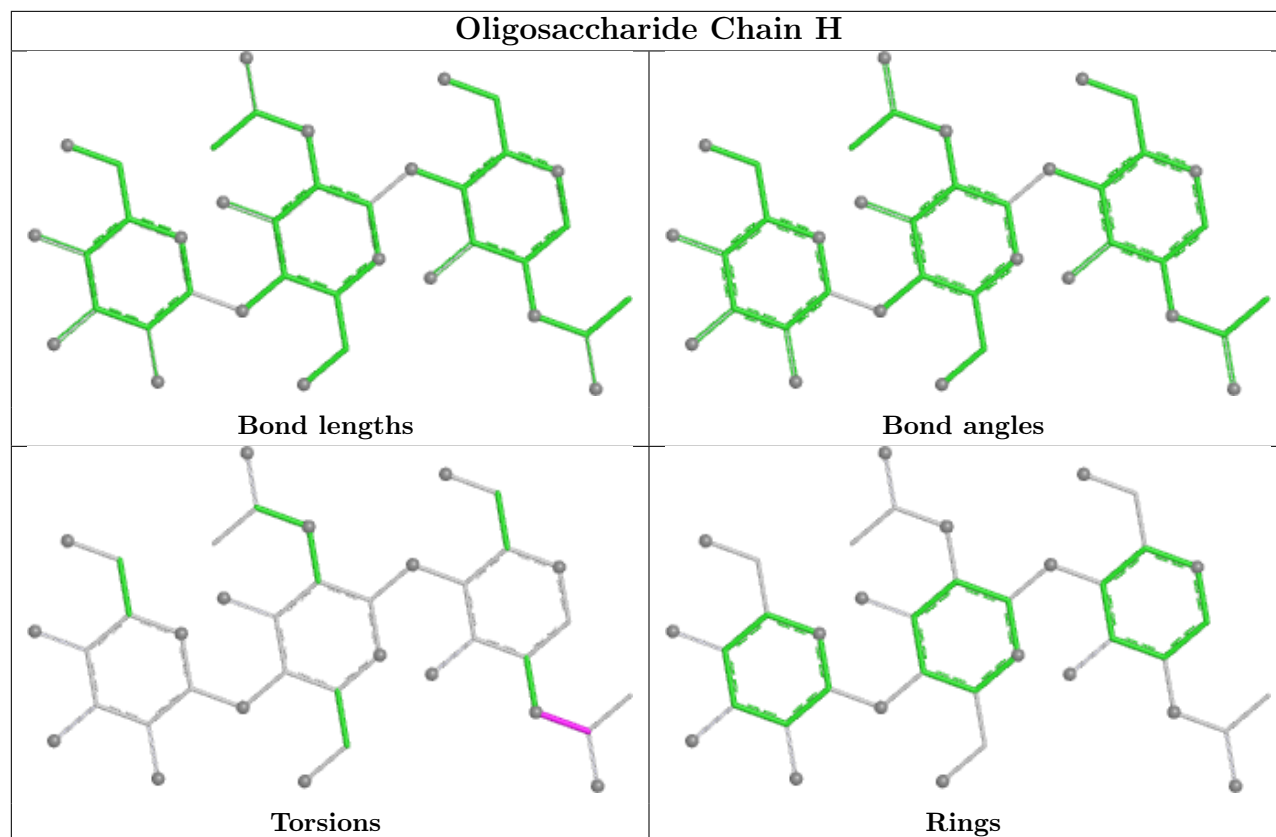


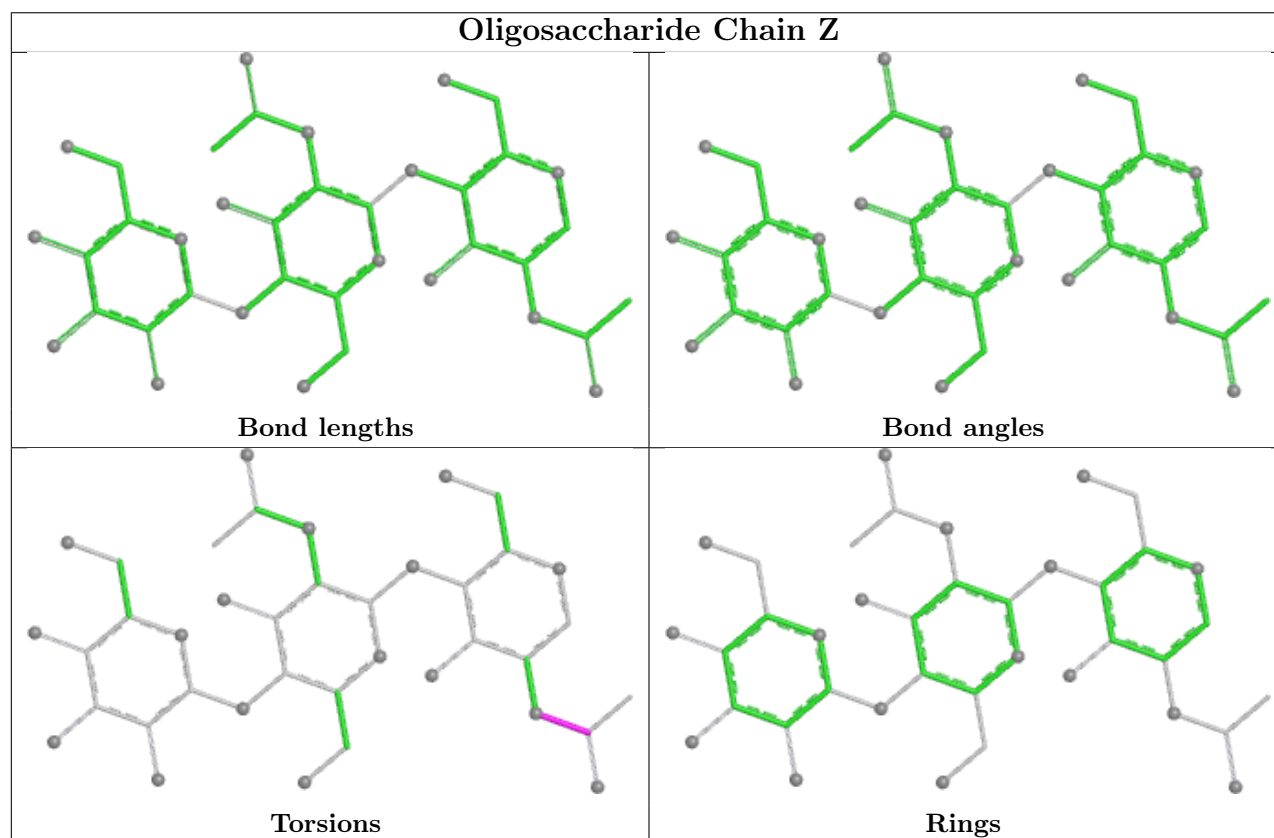
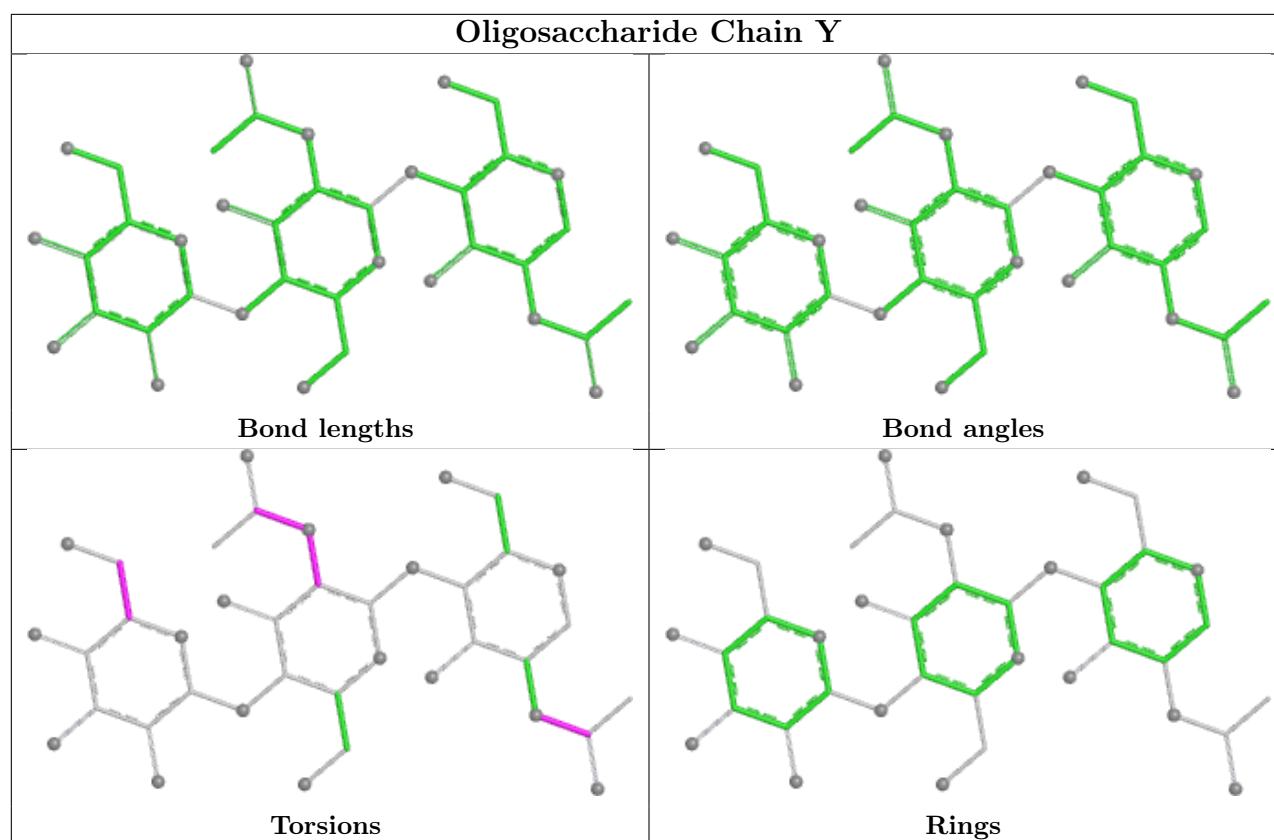


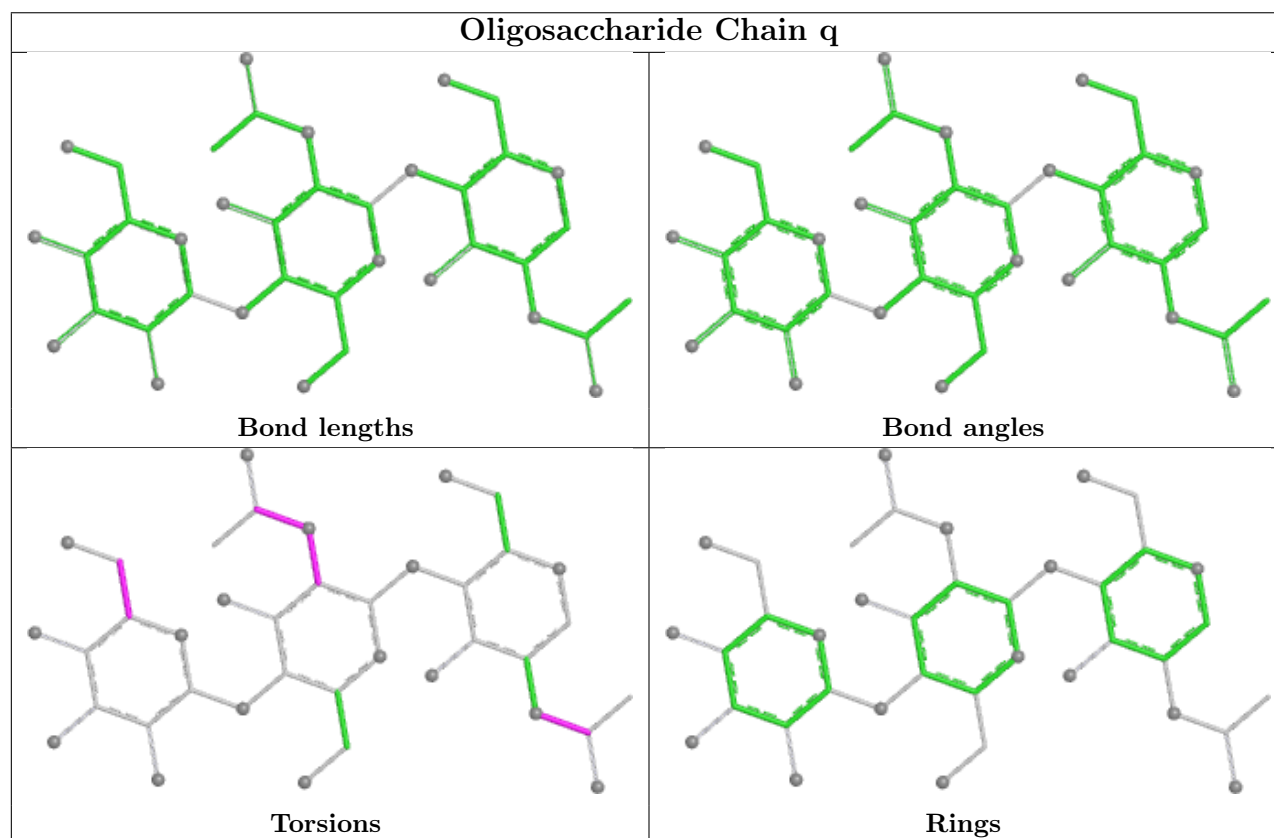
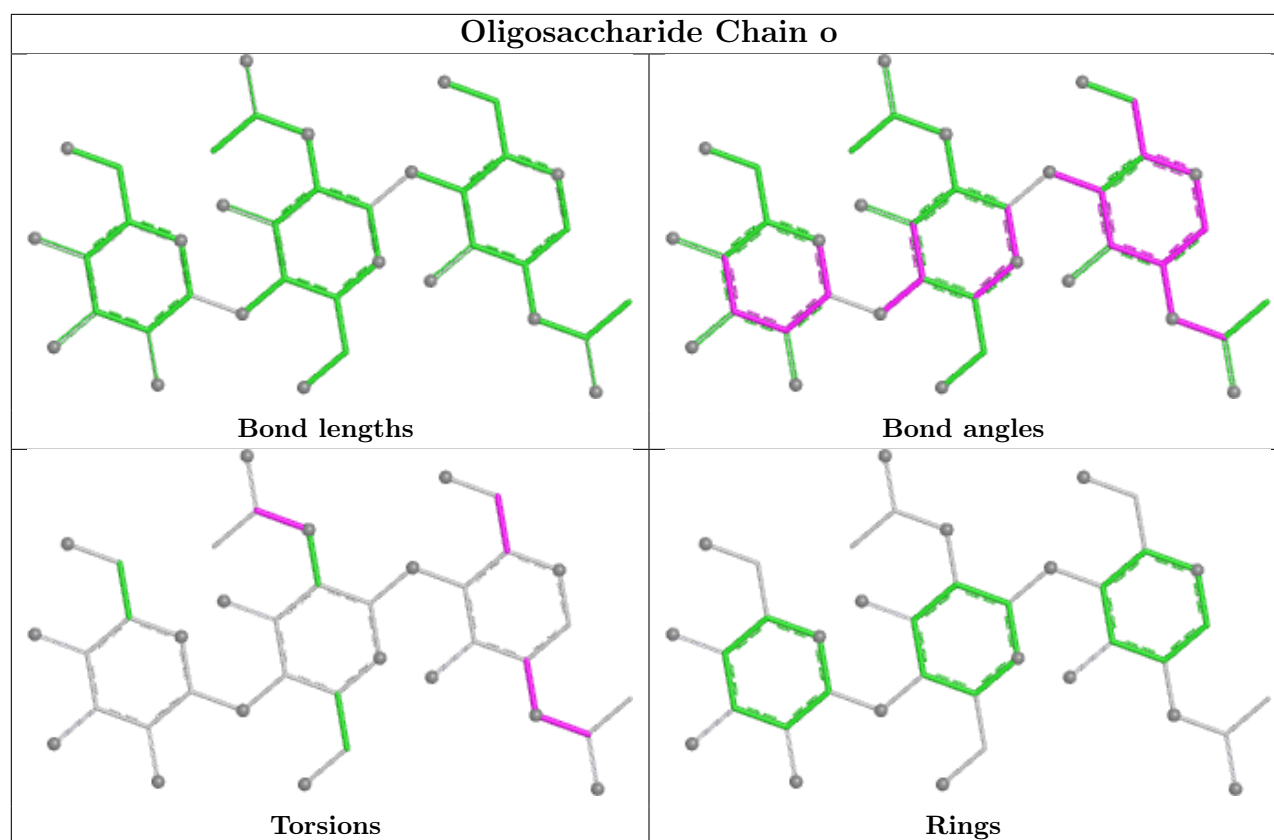


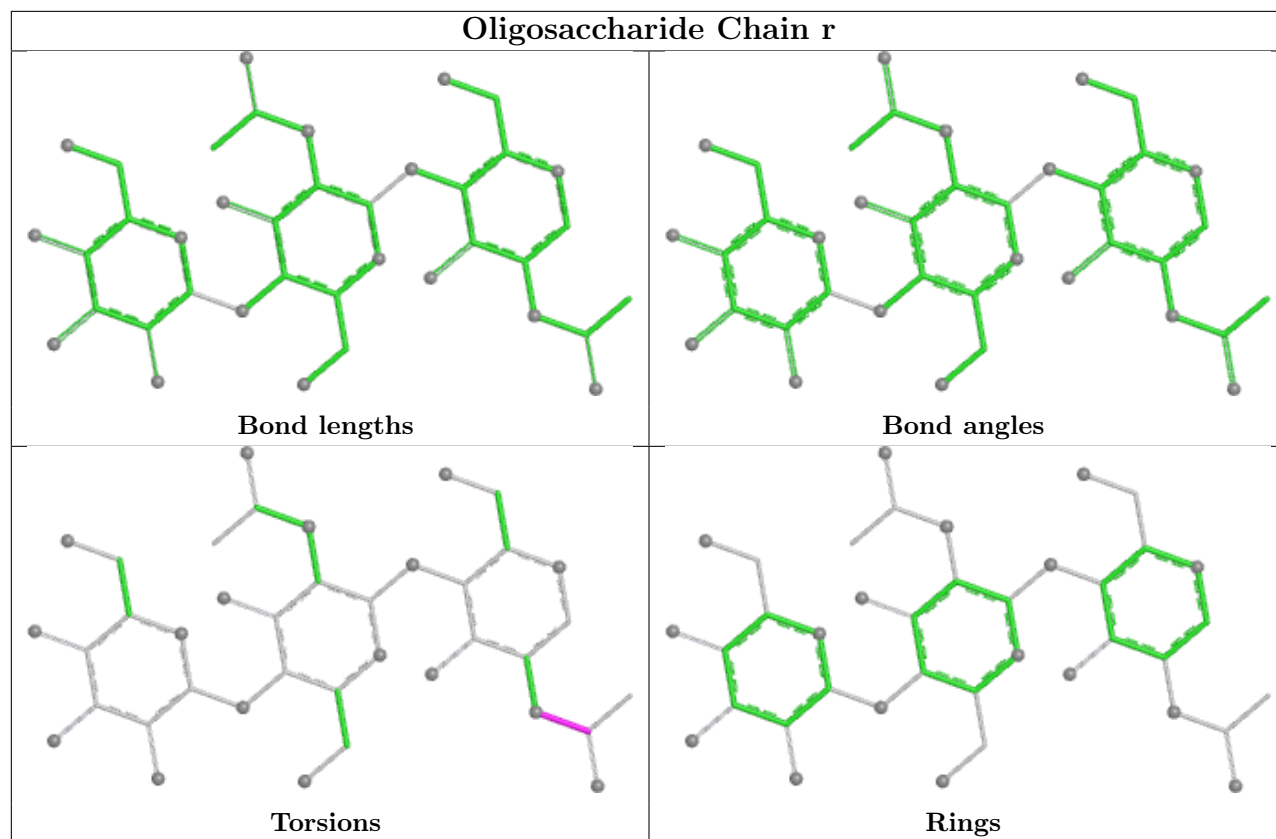


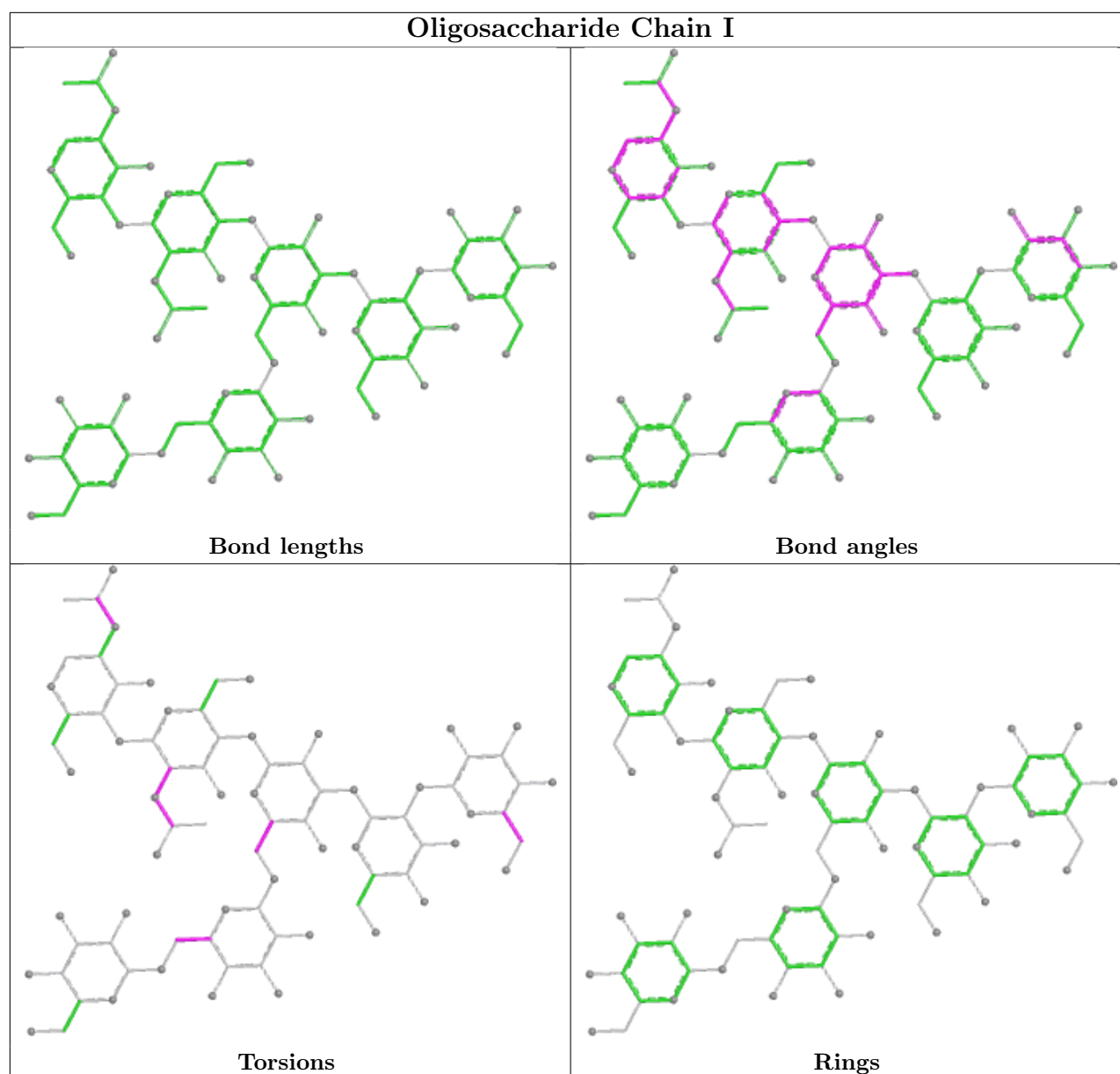


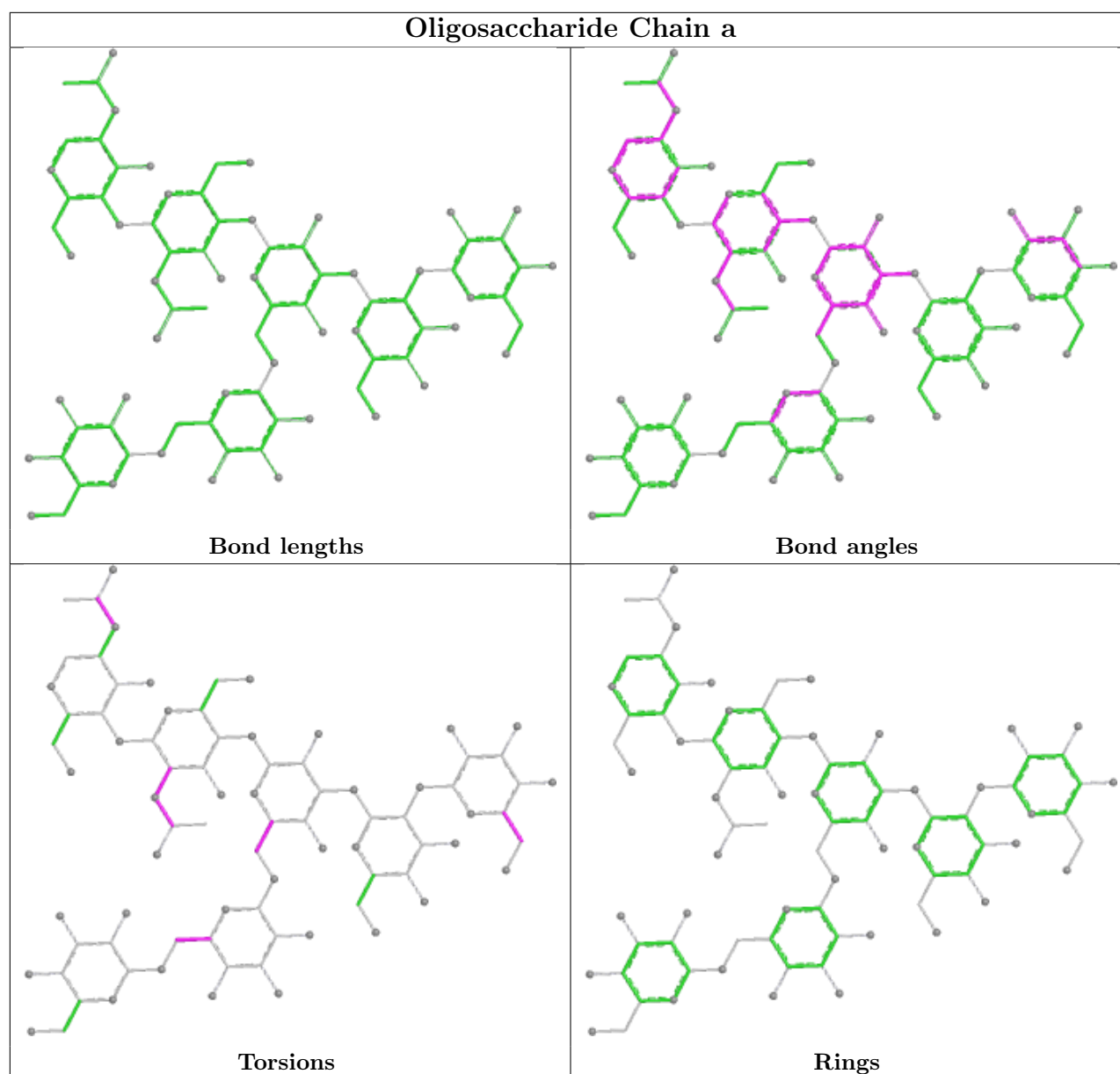


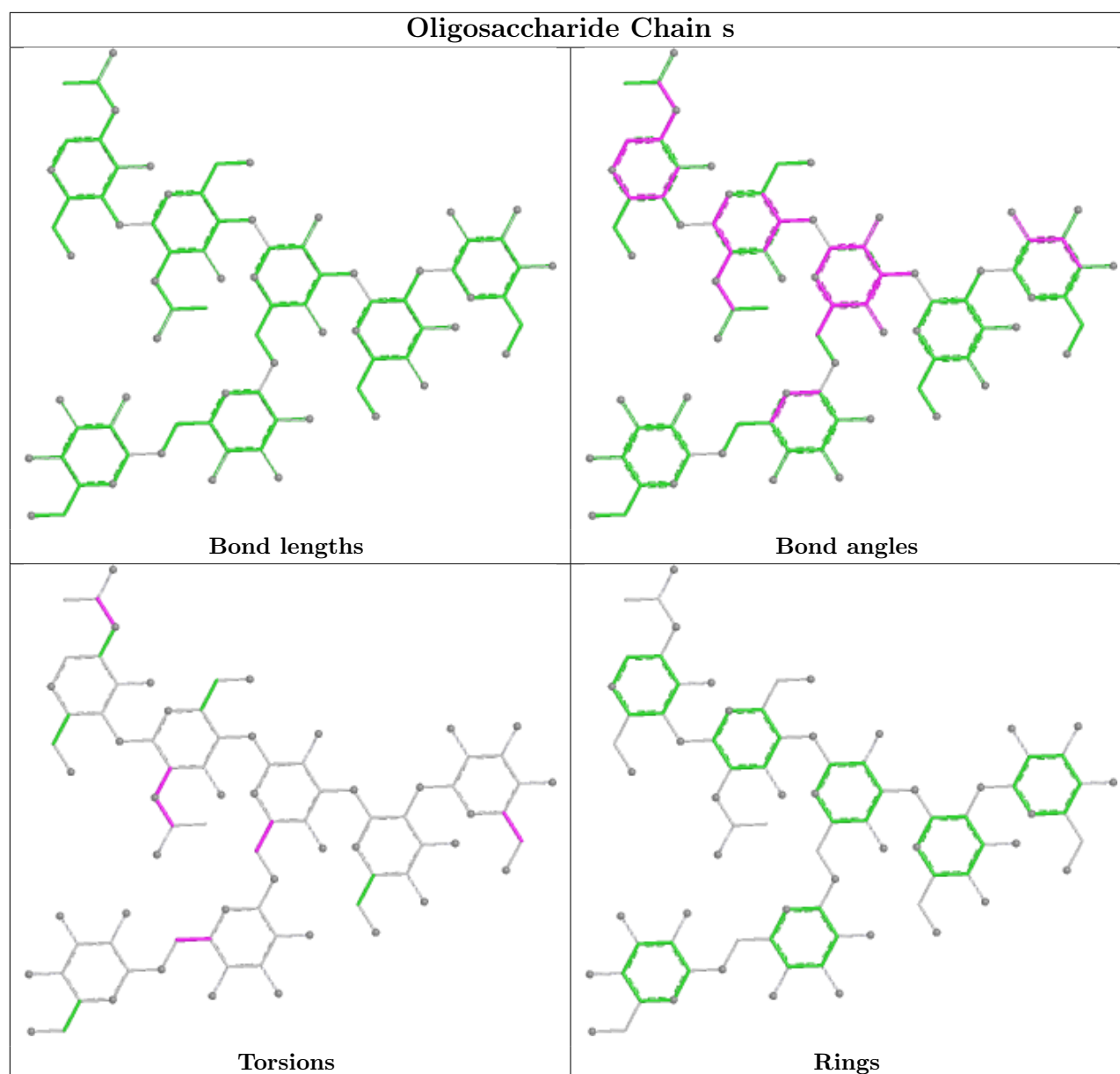


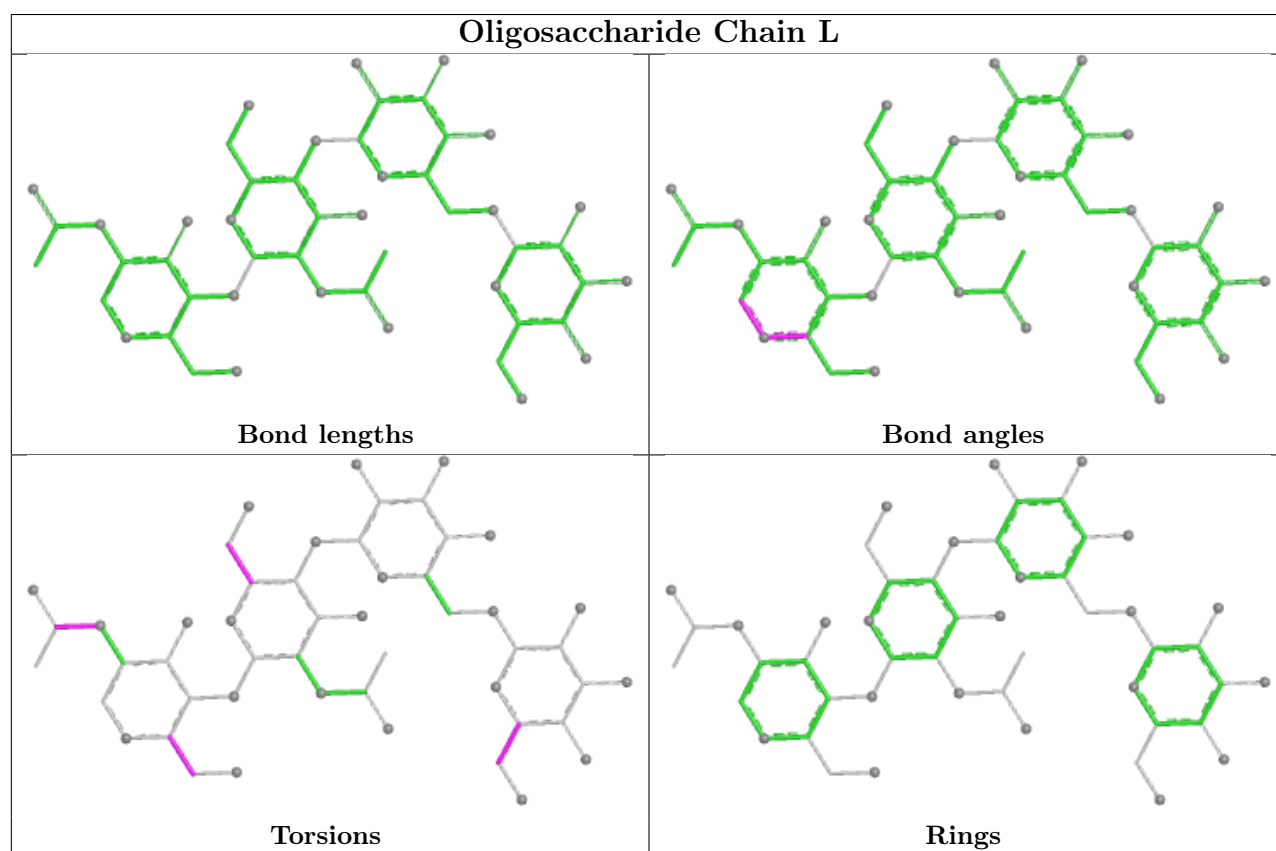
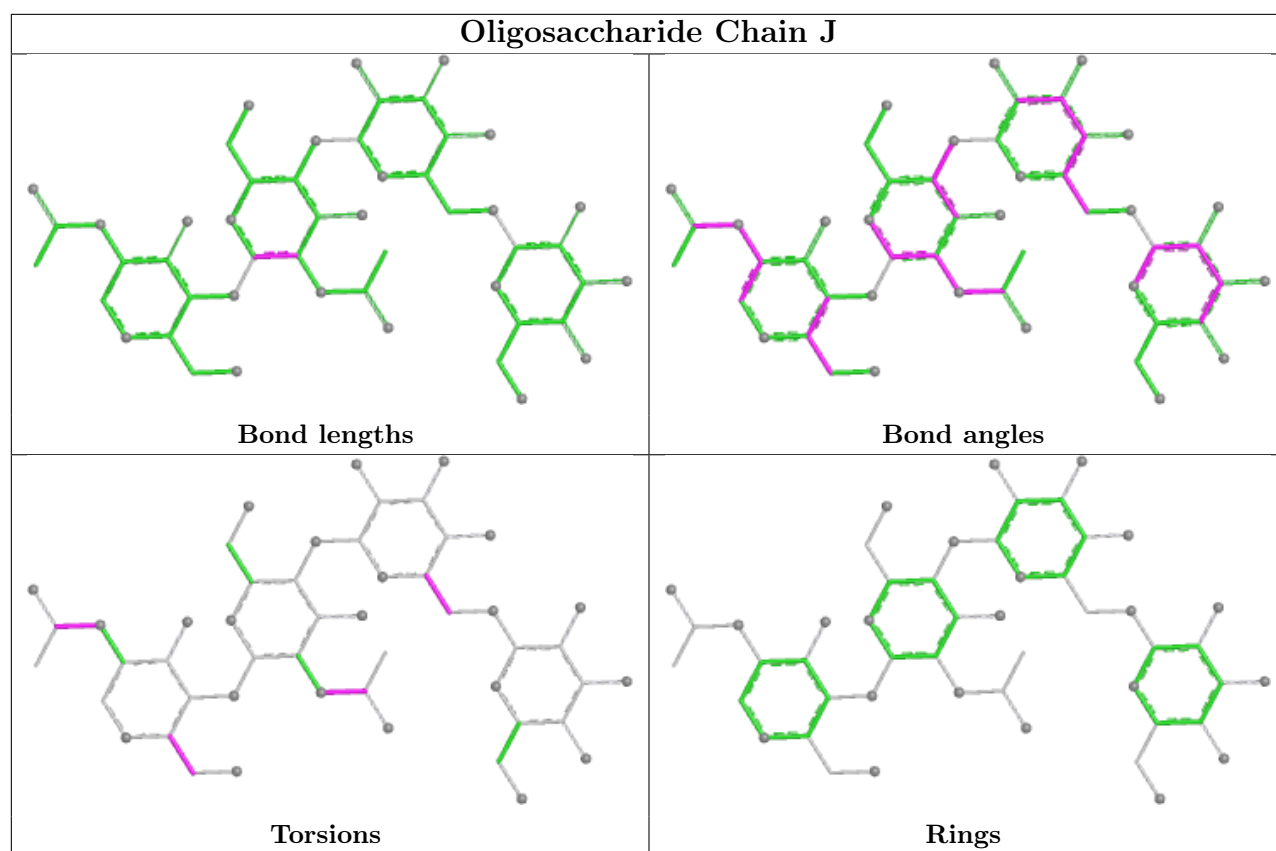


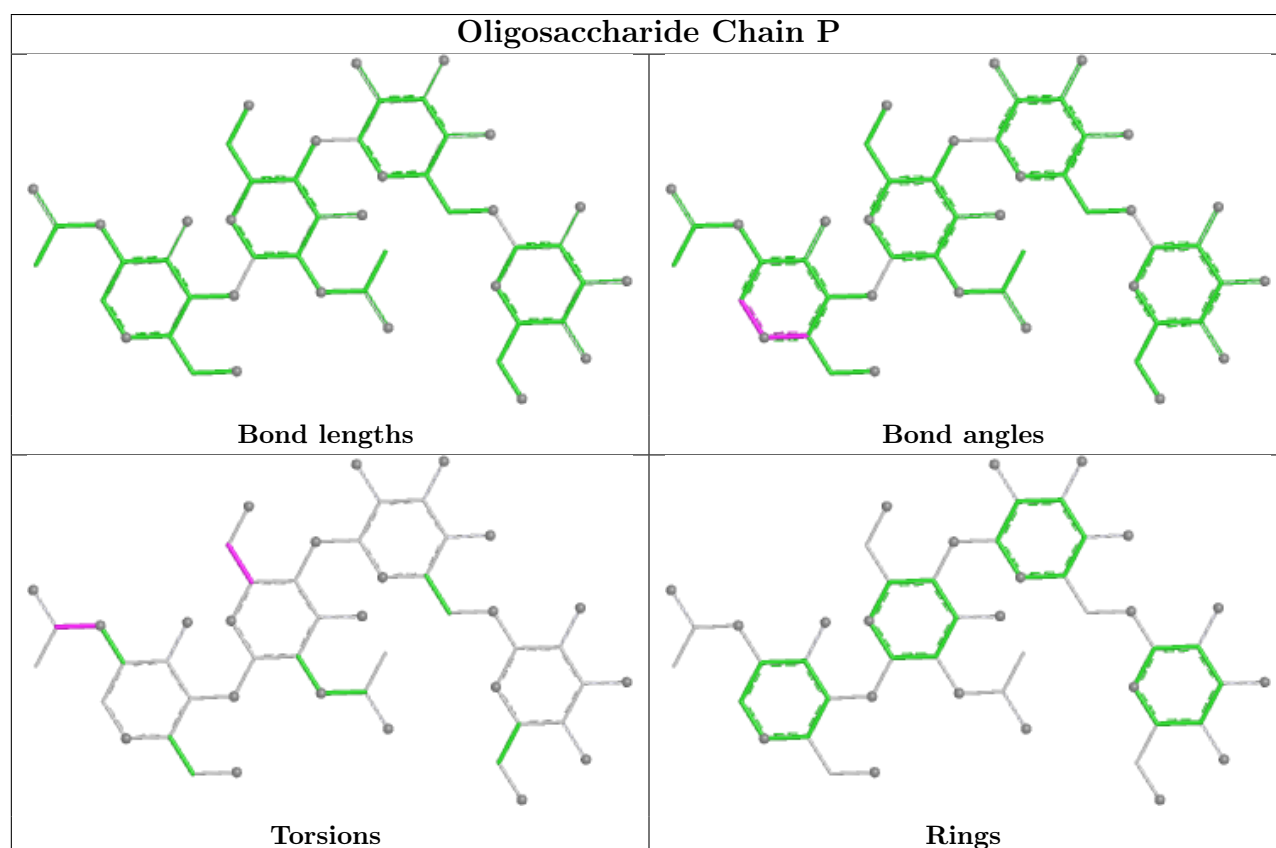
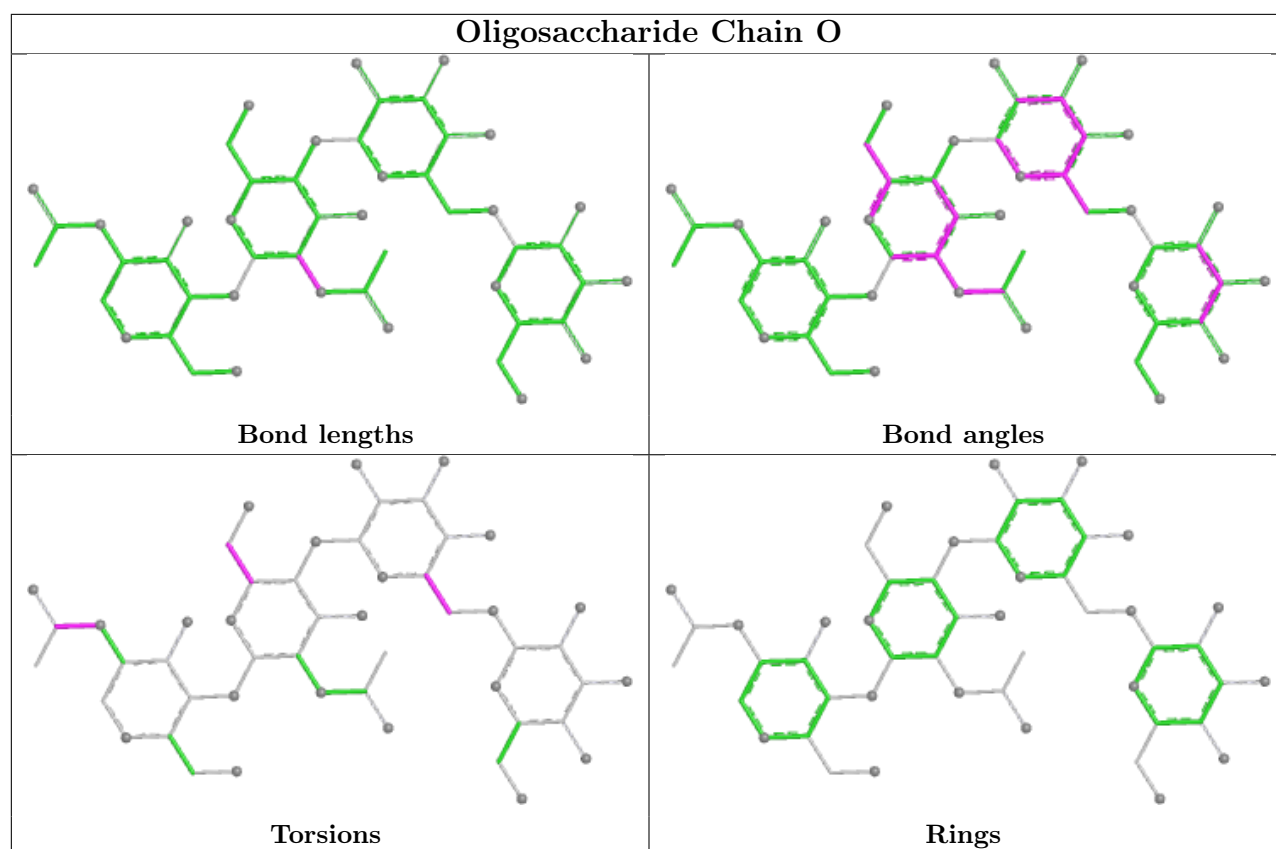


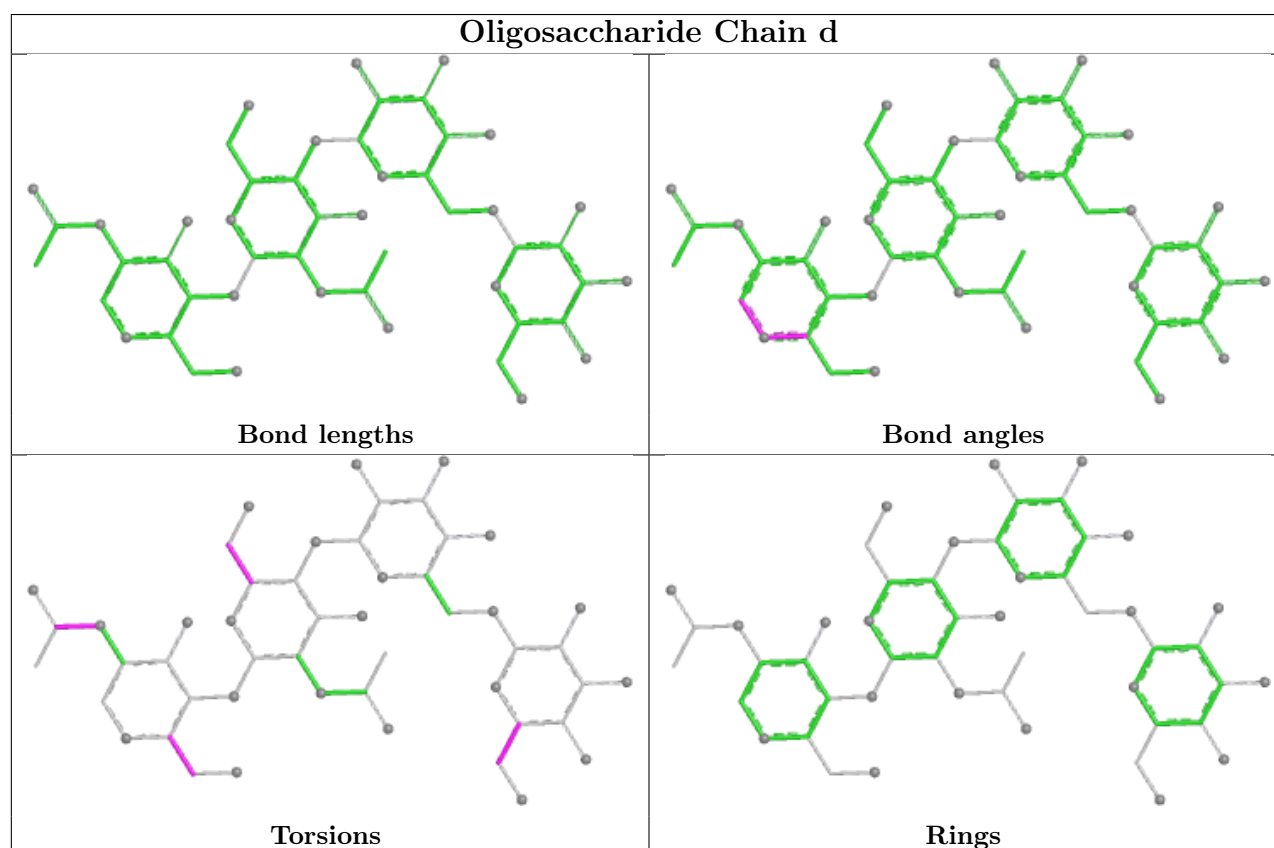
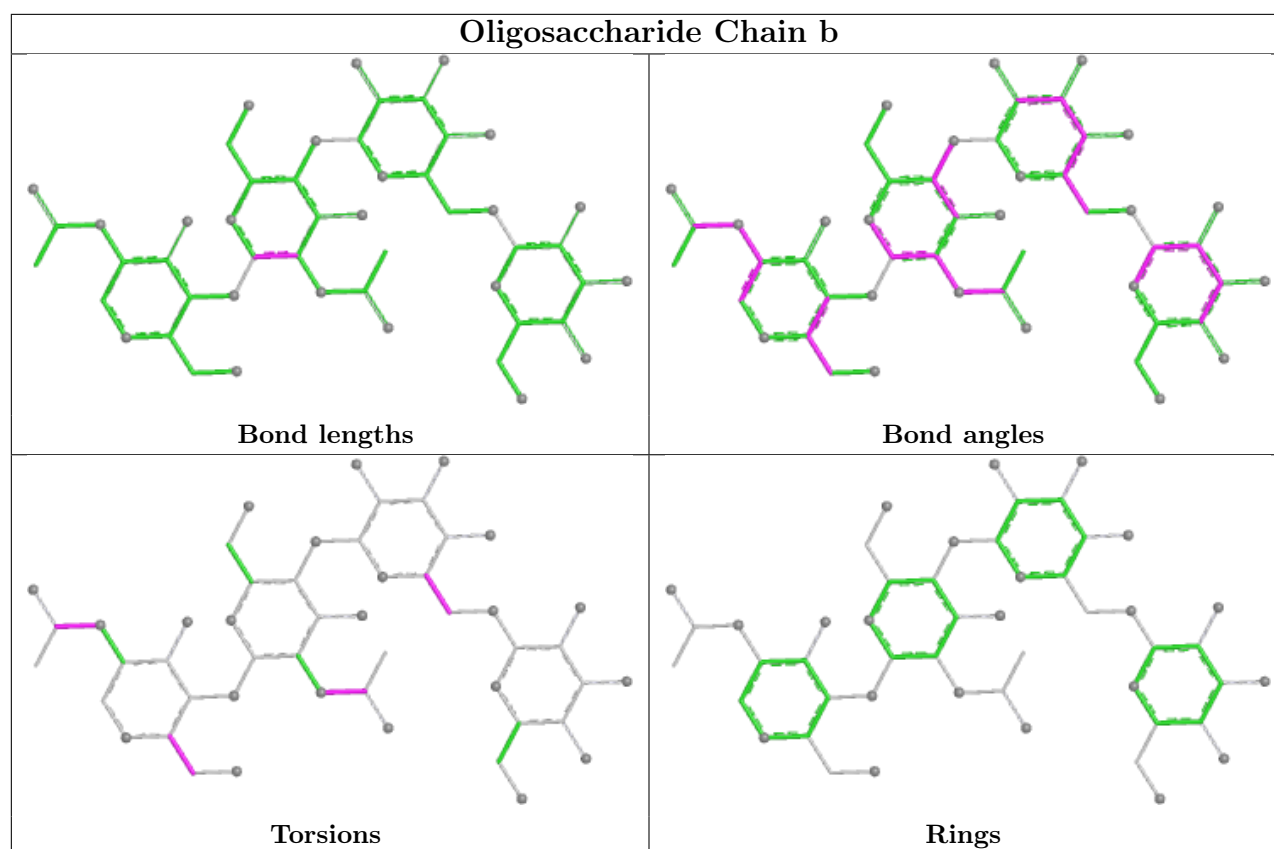


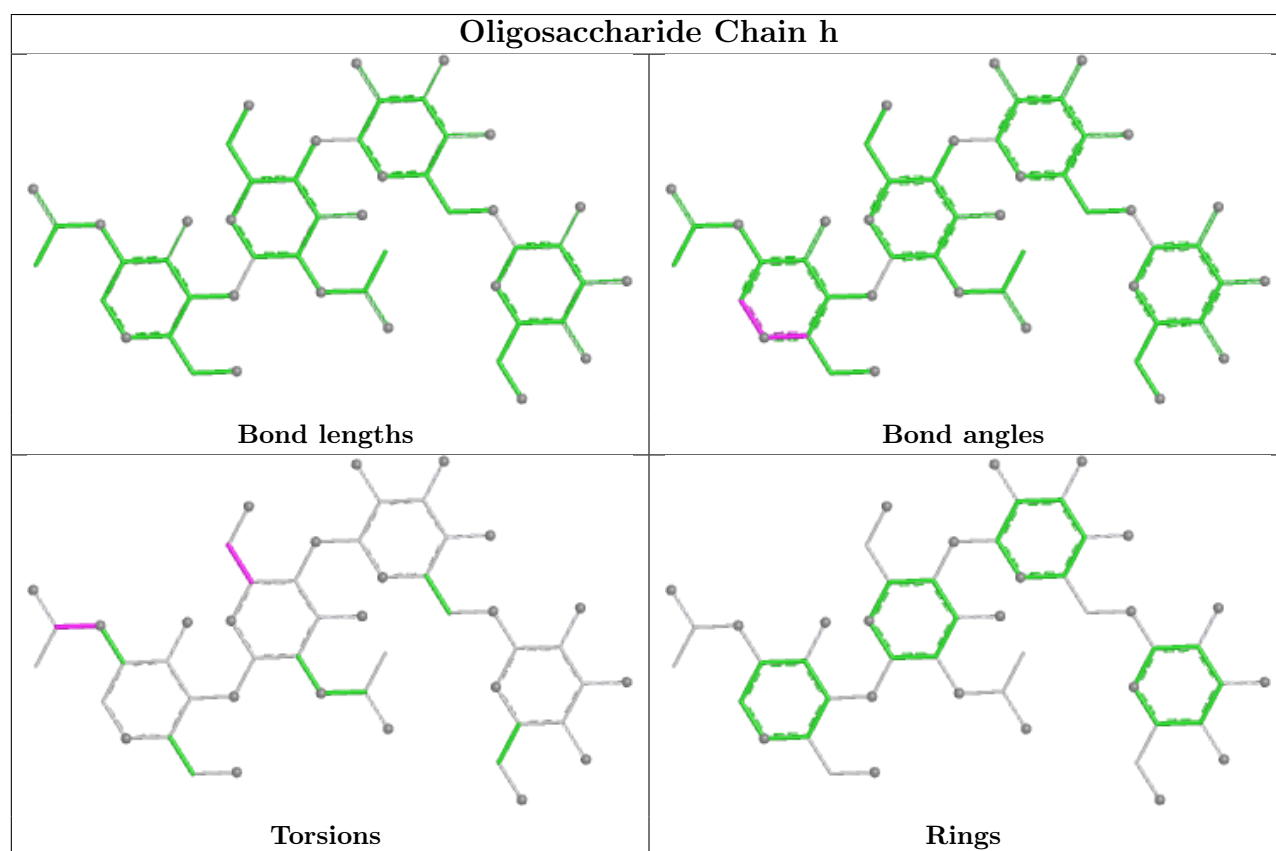
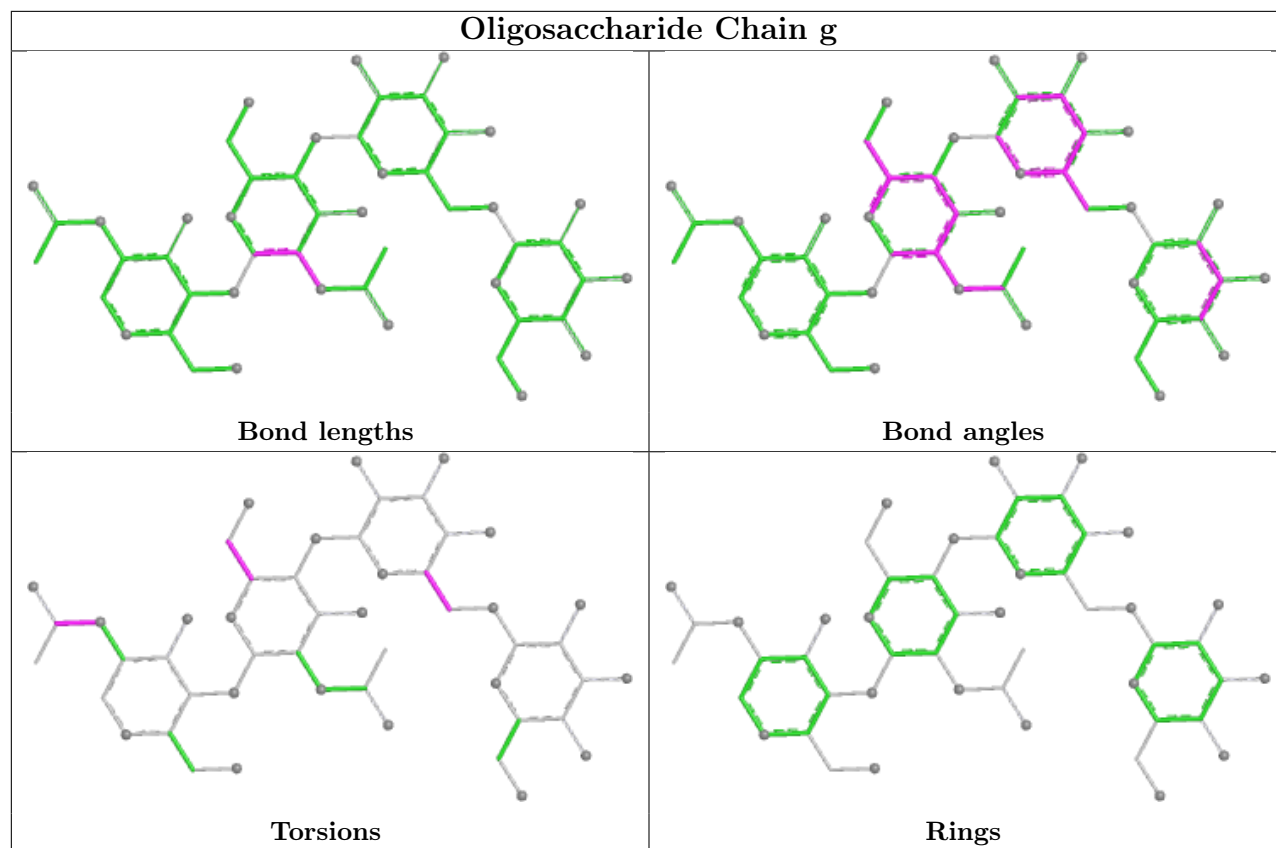


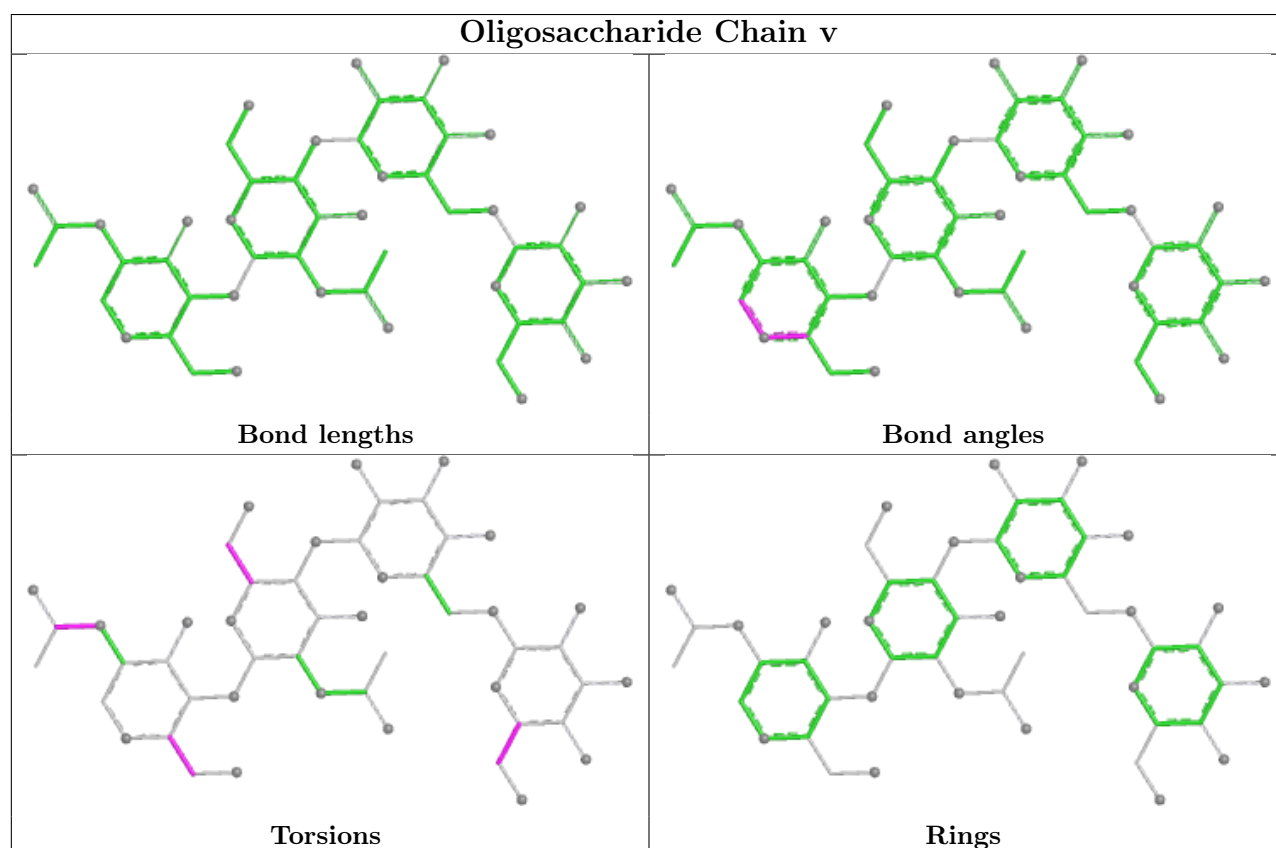
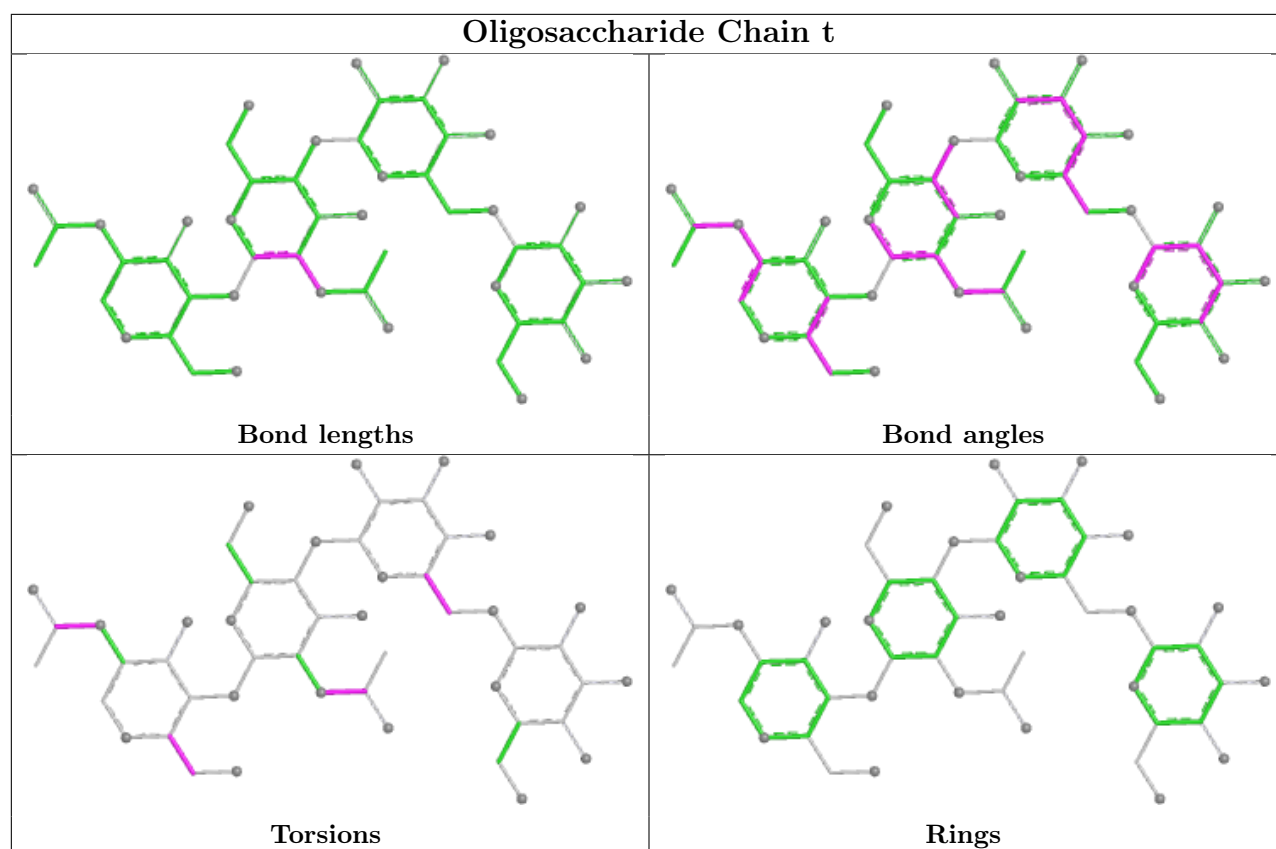


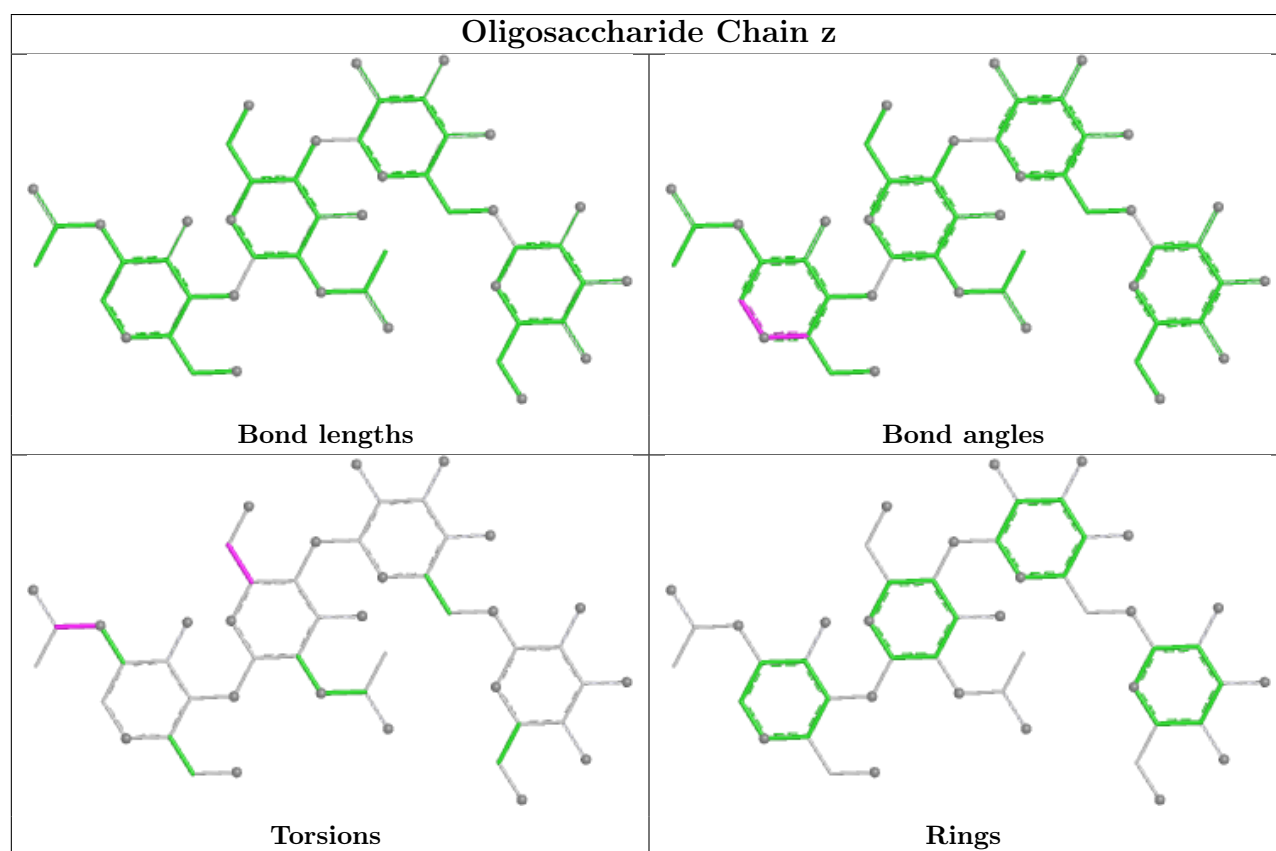
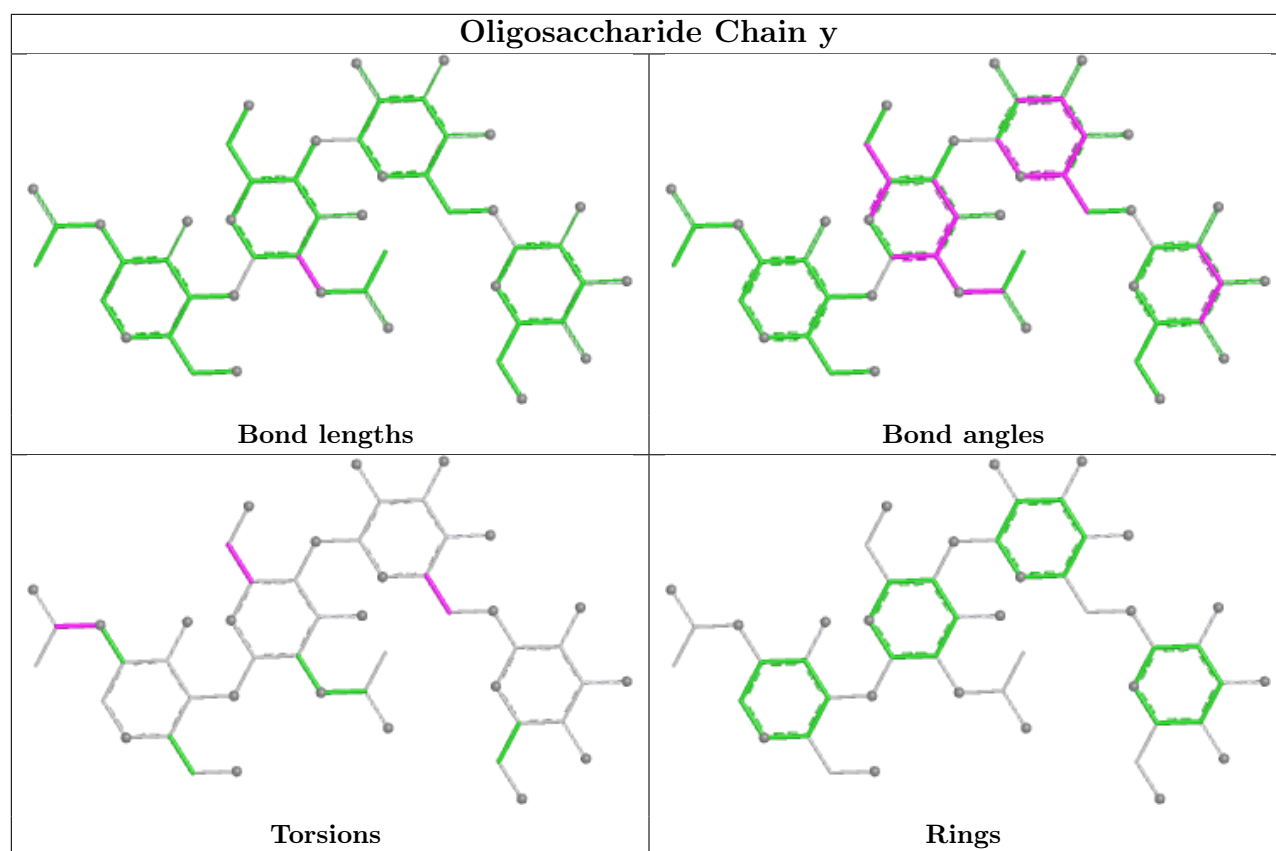


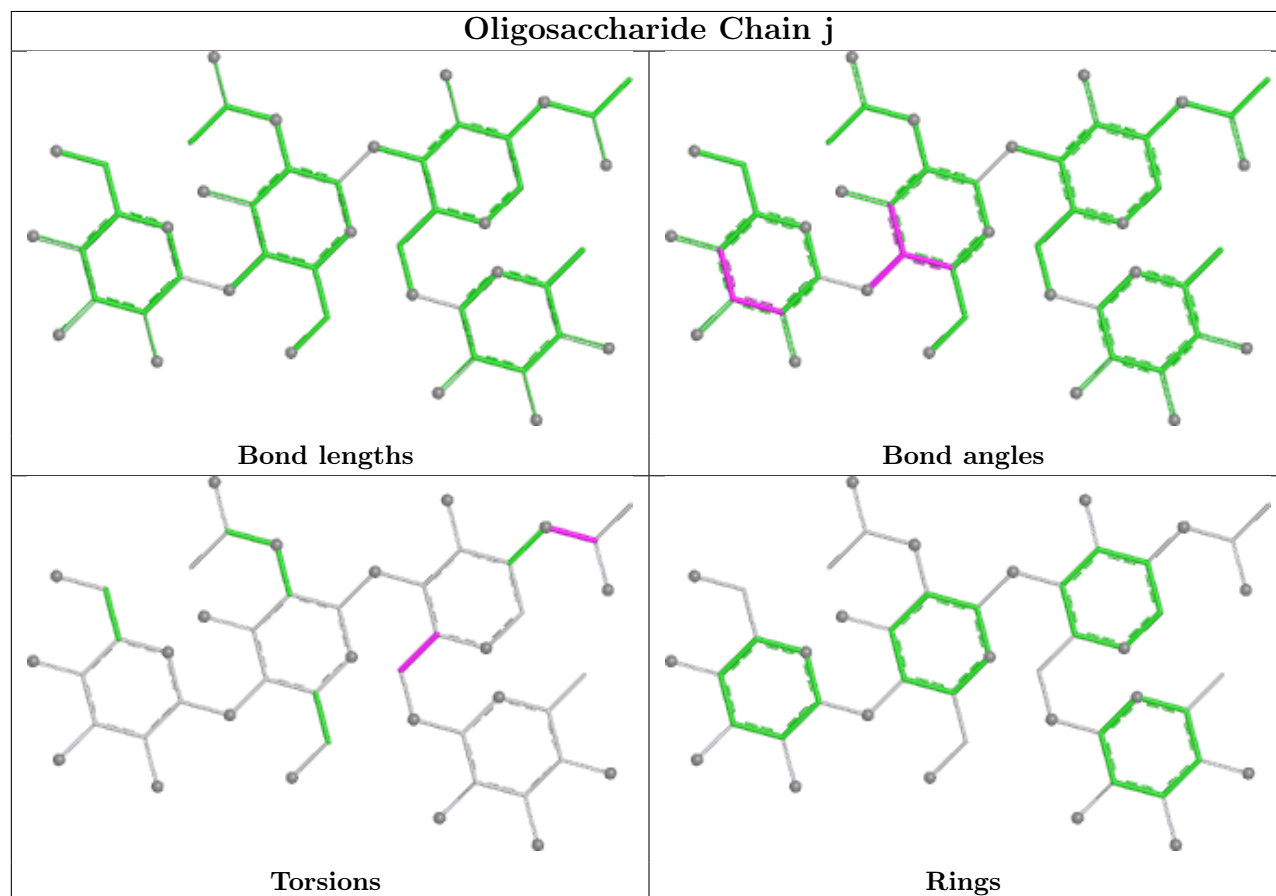
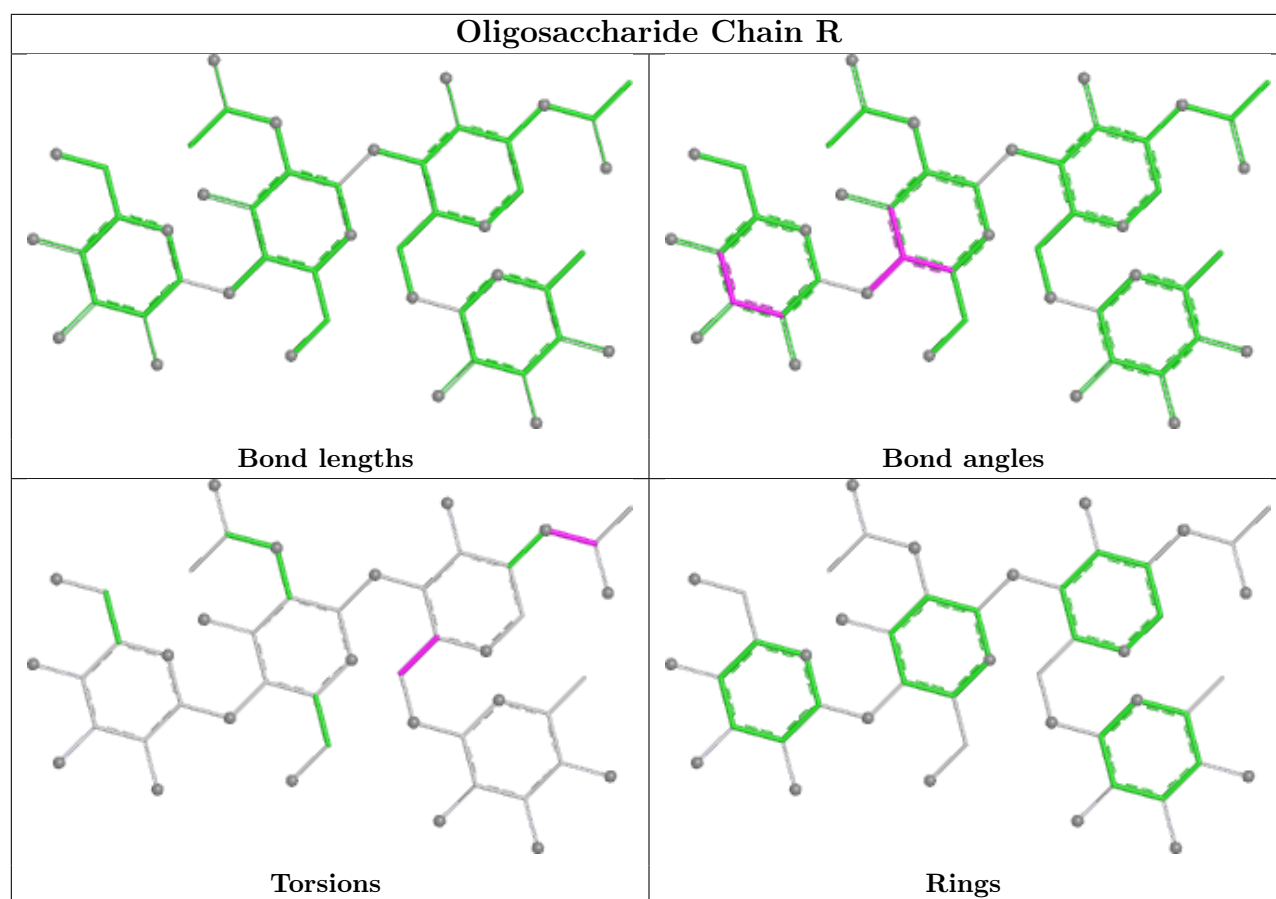


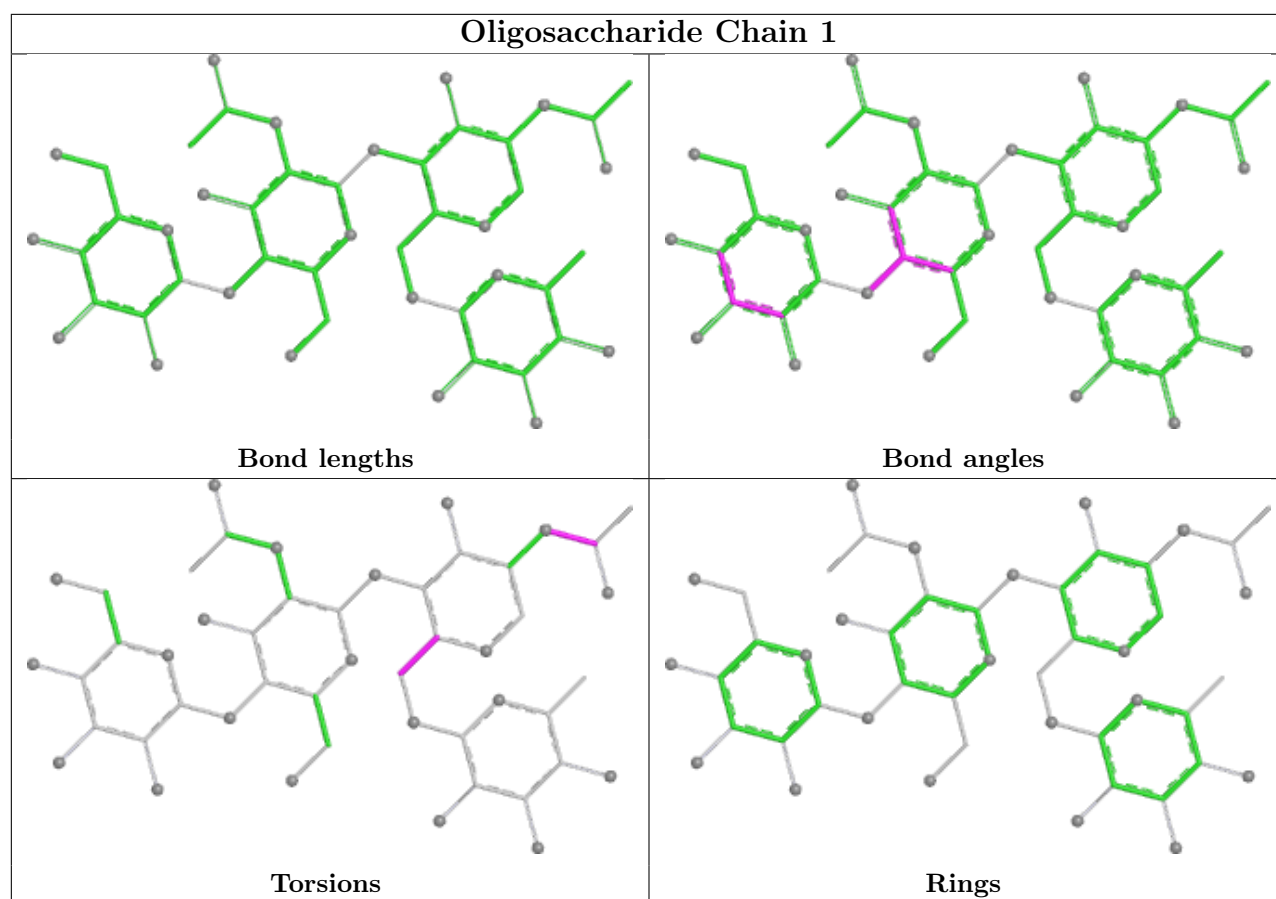


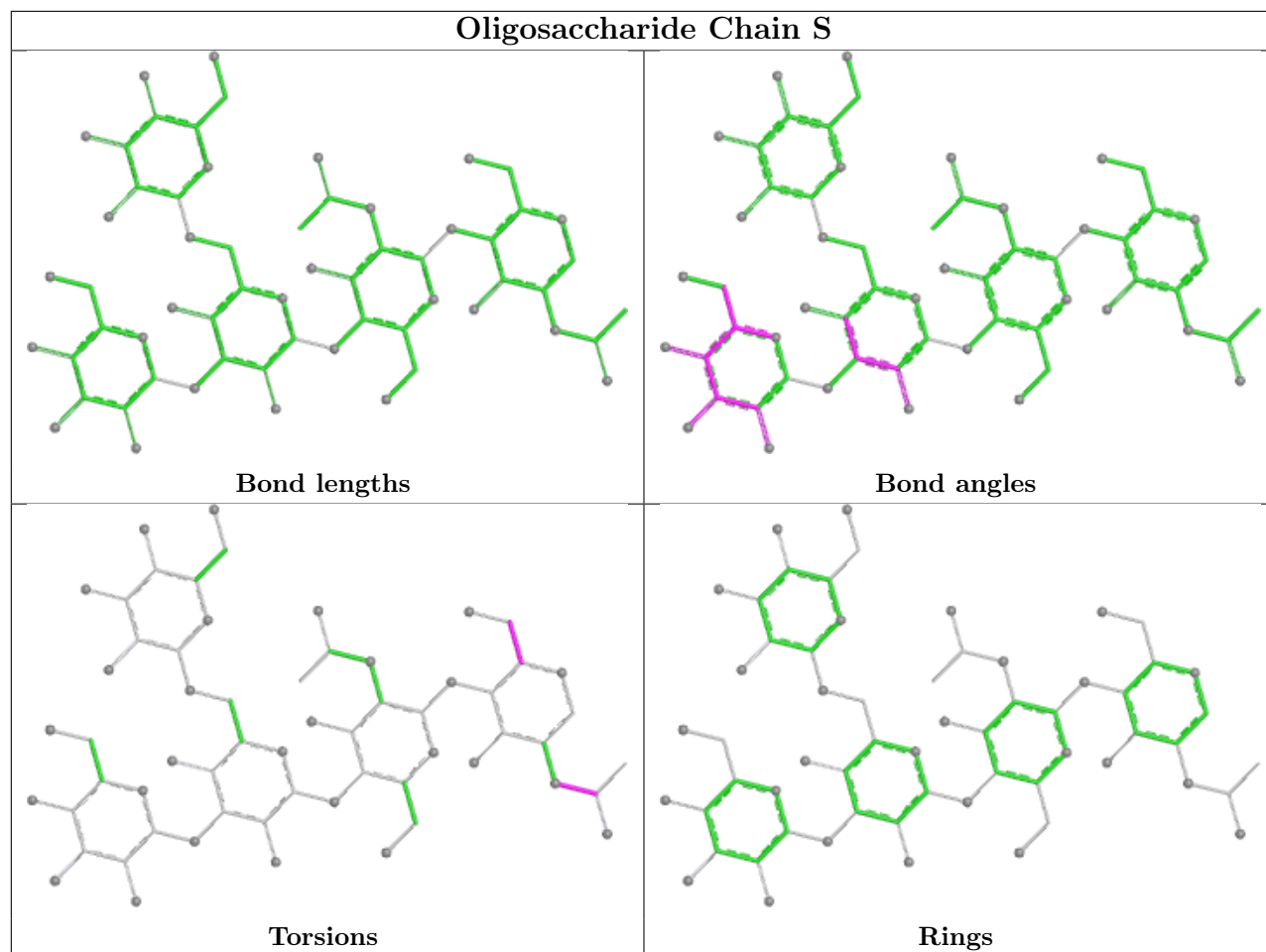


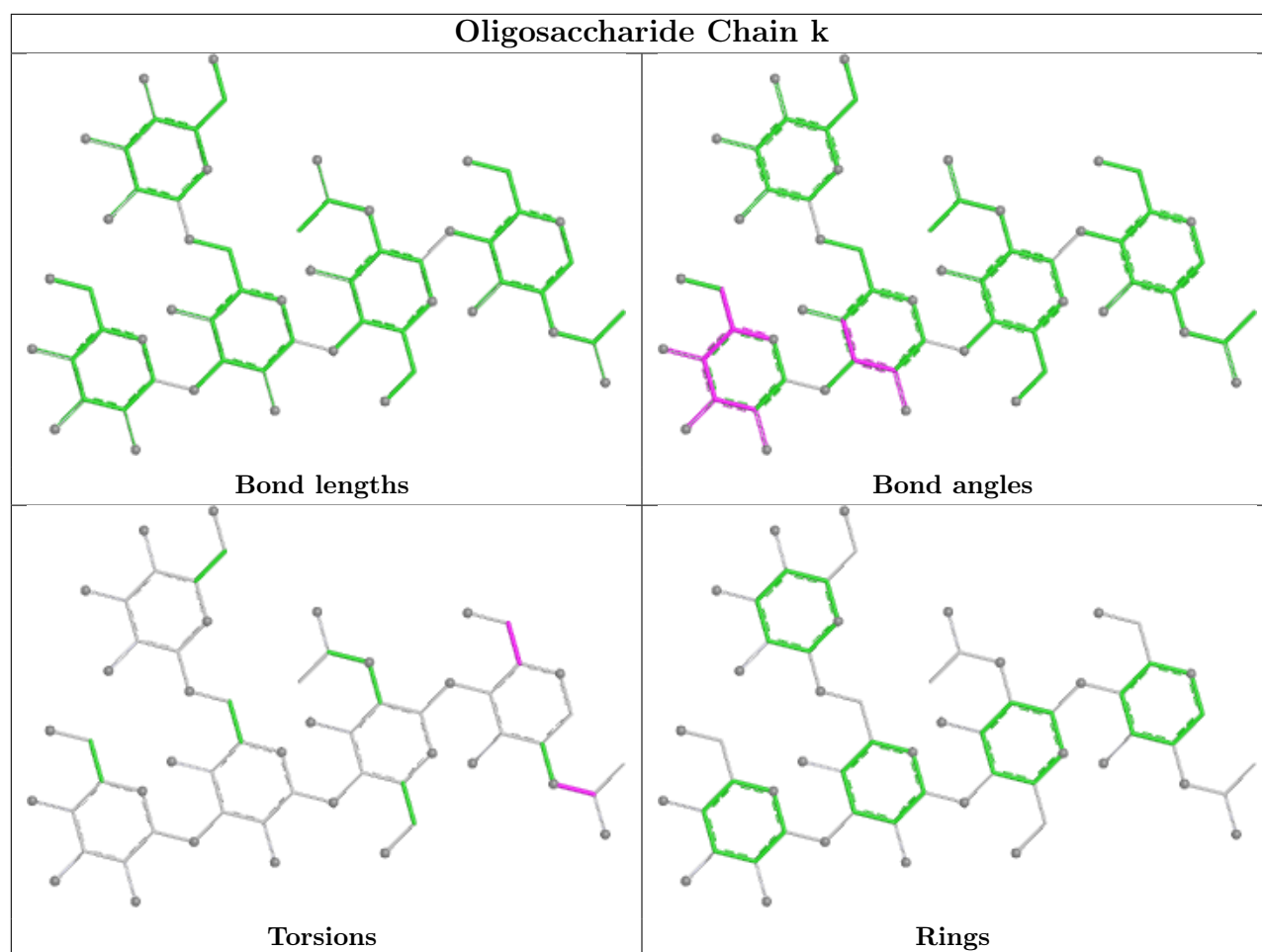


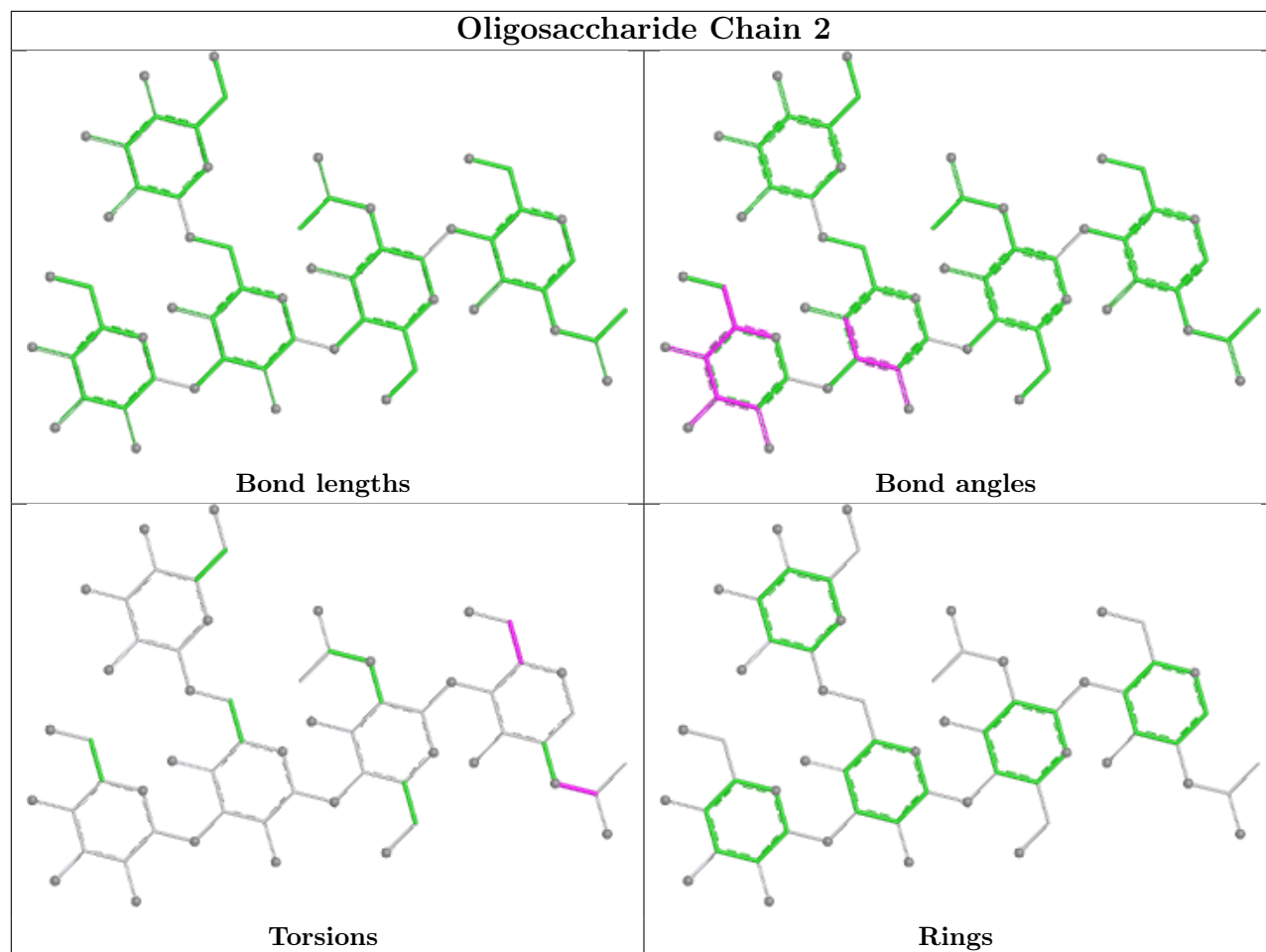


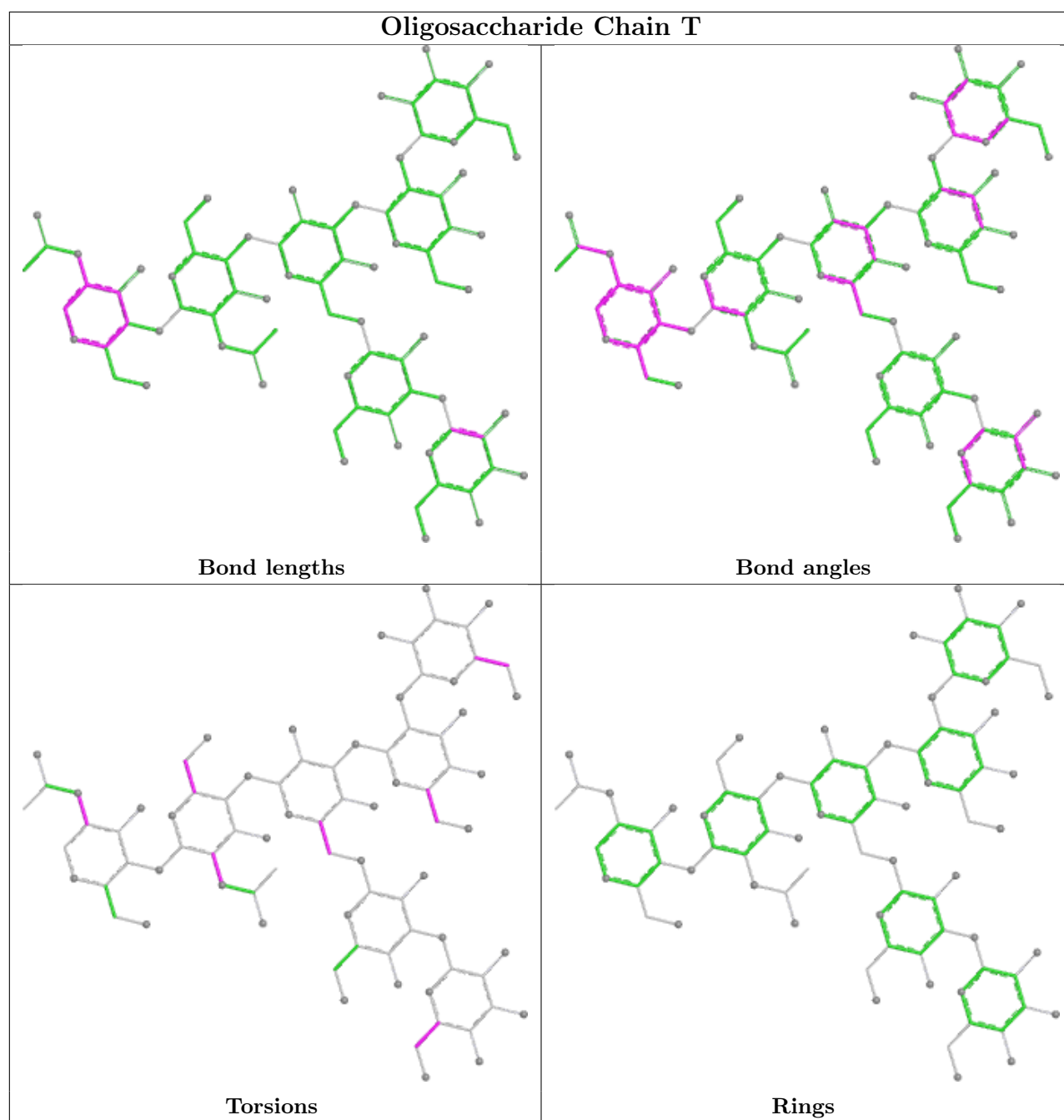


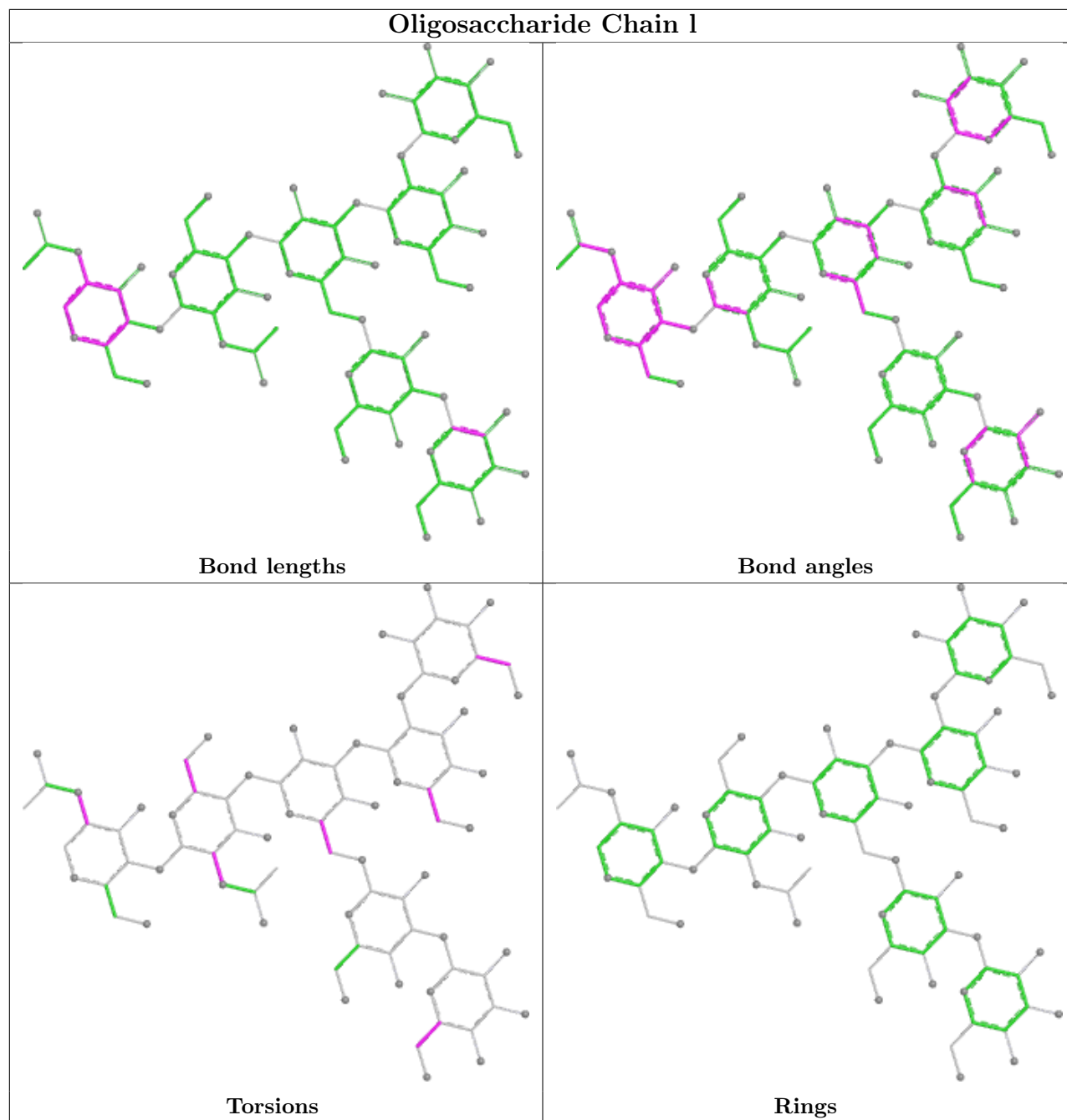


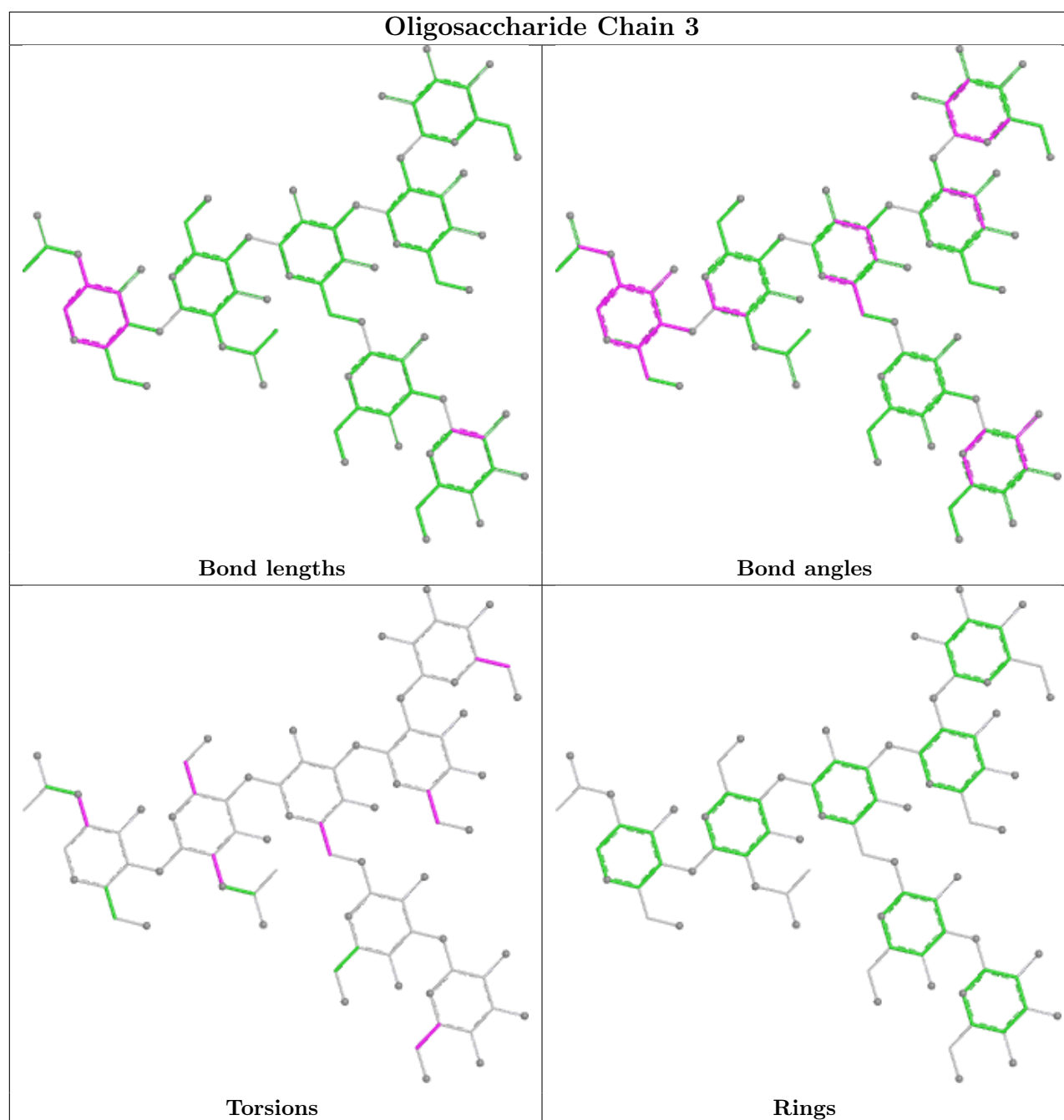












5.6 Ligand geometry [i](#)

33 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
9	NAG	B	1522	1	14,14,15	0.90	1 (7%)	17,19,21	2.34	3 (17%)
9	NAG	B	1508	1	14,14,15	0.33	0	17,19,21	0.39	0
10	MAN	C	1529	-	11,11,12	0.43	0	15,15,17	0.92	1 (6%)
9	NAG	B	1568	1	14,14,15	0.30	0	17,19,21	0.58	0
9	NAG	C	1570	1	14,14,15	0.20	0	17,19,21	0.65	0
10	MAN	B	1528	-	11,11,12	0.42	0	15,15,17	0.92	1 (6%)
9	NAG	C	1542	1	14,14,15	0.38	0	17,19,21	0.77	1 (5%)
9	NAG	A	1522	1	14,14,15	0.90	1 (7%)	17,19,21	2.34	3 (17%)
10	MAN	A	1528	-	11,11,12	0.44	0	15,15,17	0.92	1 (6%)
9	NAG	B	1535	1	14,14,15	0.28	0	17,19,21	0.72	0
9	NAG	C	1536	1	14,14,15	0.28	0	17,19,21	0.72	0
9	NAG	A	1536	1	14,14,15	0.35	0	17,19,21	0.51	0
9	NAG	C	1523	1	14,14,15	0.90	1 (7%)	17,19,21	2.34	3 (17%)
9	NAG	C	1573	1	14,14,15	0.41	0	17,19,21	0.61	0
9	NAG	A	1523	1	14,14,15	0.21	0	17,19,21	0.49	0
9	NAG	C	1524	1	14,14,15	0.22	0	17,19,21	0.49	0
9	NAG	B	1541	1	14,14,15	0.38	0	17,19,21	0.77	1 (5%)
9	NAG	A	1568	1	14,14,15	0.30	0	17,19,21	0.58	0
9	NAG	B	1569	1	14,14,15	0.21	0	17,19,21	0.65	0
9	NAG	A	1508	1	14,14,15	0.33	0	17,19,21	0.39	0
9	NAG	A	1541	1	14,14,15	0.38	0	17,19,21	0.77	1 (5%)
9	NAG	A	1569	1	14,14,15	0.20	0	17,19,21	0.65	0
9	NAG	C	1509	1	14,14,15	0.33	0	17,19,21	0.39	0
9	NAG	B	1536	1	14,14,15	0.36	0	17,19,21	0.51	0
10	MAN	C	1501	-	11,11,12	0.46	0	15,15,17	1.23	1 (6%)
9	NAG	A	1535	1	14,14,15	0.28	0	17,19,21	0.72	0
10	MAN	B	1573	-	11,11,12	0.45	0	15,15,17	1.23	1 (6%)
9	NAG	B	1523	1	14,14,15	0.21	0	17,19,21	0.49	0
9	NAG	A	1572	1	14,14,15	0.42	0	17,19,21	0.60	0
9	NAG	C	1569	1	14,14,15	0.30	0	17,19,21	0.58	0
9	NAG	B	1572	1	14,14,15	0.42	0	17,19,21	0.60	0
9	NAG	C	1537	1	14,14,15	0.36	0	17,19,21	0.51	0
10	MAN	A	1573	-	11,11,12	0.45	0	15,15,17	1.23	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
9	NAG	B	1522	1	-	6/6/23/26	0/1/1/1
9	NAG	B	1508	1	-	2/6/23/26	0/1/1/1
10	MAN	C	1529	-	-	0/2/19/22	0/1/1/1
9	NAG	B	1568	1	-	4/6/23/26	0/1/1/1
9	NAG	C	1570	1	-	4/6/23/26	0/1/1/1
10	MAN	B	1528	-	-	0/2/19/22	0/1/1/1
9	NAG	C	1542	1	-	4/6/23/26	0/1/1/1
9	NAG	A	1522	1	-	6/6/23/26	0/1/1/1
10	MAN	A	1528	-	-	0/2/19/22	0/1/1/1
9	NAG	B	1535	1	-	2/6/23/26	0/1/1/1
9	NAG	C	1536	1	-	2/6/23/26	0/1/1/1
9	NAG	A	1536	1	-	0/6/23/26	0/1/1/1
9	NAG	C	1523	1	-	6/6/23/26	0/1/1/1
9	NAG	C	1573	1	-	4/6/23/26	0/1/1/1
9	NAG	A	1523	1	-	2/6/23/26	0/1/1/1
9	NAG	C	1524	1	-	2/6/23/26	0/1/1/1
9	NAG	B	1541	1	-	4/6/23/26	0/1/1/1
9	NAG	A	1568	1	-	4/6/23/26	0/1/1/1
9	NAG	B	1569	1	-	4/6/23/26	0/1/1/1
9	NAG	A	1508	1	-	2/6/23/26	0/1/1/1
9	NAG	A	1541	1	-	4/6/23/26	0/1/1/1
9	NAG	A	1569	1	-	4/6/23/26	0/1/1/1
9	NAG	C	1509	1	-	2/6/23/26	0/1/1/1
9	NAG	B	1536	1	-	0/6/23/26	0/1/1/1
10	MAN	C	1501	-	-	0/2/19/22	0/1/1/1
9	NAG	A	1535	1	-	2/6/23/26	0/1/1/1
10	MAN	B	1573	-	-	0/2/19/22	0/1/1/1
9	NAG	B	1523	1	-	2/6/23/26	0/1/1/1
9	NAG	A	1572	1	-	4/6/23/26	0/1/1/1
9	NAG	C	1569	1	-	4/6/23/26	0/1/1/1
9	NAG	B	1572	1	-	4/6/23/26	0/1/1/1
9	NAG	C	1537	1	-	0/6/23/26	0/1/1/1
10	MAN	A	1573	-	-	0/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	1522	NAG	C1-C2	2.60	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	1522	NAG	C1-C2	2.60	1.55	1.52
9	C	1523	NAG	C1-C2	2.60	1.55	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	1522	NAG	C2-N2-C7	8.36	134.10	122.90
9	A	1522	NAG	C2-N2-C7	8.35	134.09	122.90
9	C	1523	NAG	C2-N2-C7	8.34	134.08	122.90
9	B	1522	NAG	C1-C2-N2	3.82	116.46	110.43
9	A	1522	NAG	C1-C2-N2	3.82	116.45	110.43
9	C	1523	NAG	C1-C2-N2	3.81	116.43	110.43
9	A	1541	NAG	C1-O5-C5	2.43	115.45	112.19
9	B	1541	NAG	C1-O5-C5	2.42	115.44	112.19
9	C	1542	NAG	C1-O5-C5	2.42	115.42	112.19
10	B	1528	MAN	C2-C3-C4	-2.36	106.70	110.86
10	C	1529	MAN	C2-C3-C4	-2.36	106.72	110.86
10	A	1528	MAN	C2-C3-C4	-2.35	106.73	110.86
10	A	1573	MAN	C2-C3-C4	-2.25	106.91	110.86
10	B	1573	MAN	C2-C3-C4	-2.23	106.93	110.86
10	C	1501	MAN	C2-C3-C4	-2.23	106.94	110.86
9	C	1523	NAG	C8-C7-N2	2.18	119.74	116.12
9	A	1522	NAG	C8-C7-N2	2.16	119.70	116.12
9	B	1522	NAG	C8-C7-N2	2.14	119.66	116.12

There are no chirality outliers.

All (84) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	1541	NAG	O5-C5-C6-O6
9	B	1541	NAG	O5-C5-C6-O6
9	C	1542	NAG	O5-C5-C6-O6
9	A	1508	NAG	O5-C5-C6-O6
9	B	1508	NAG	O5-C5-C6-O6
9	C	1509	NAG	O5-C5-C6-O6
9	A	1522	NAG	O5-C5-C6-O6
9	B	1522	NAG	O5-C5-C6-O6
9	C	1523	NAG	O5-C5-C6-O6
9	A	1508	NAG	C4-C5-C6-O6
9	B	1508	NAG	C4-C5-C6-O6
9	C	1509	NAG	C4-C5-C6-O6
9	A	1569	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
9	B	1569	NAG	O5-C5-C6-O6
9	C	1570	NAG	O5-C5-C6-O6
9	A	1522	NAG	C4-C5-C6-O6
9	A	1541	NAG	C4-C5-C6-O6
9	B	1522	NAG	C4-C5-C6-O6
9	B	1541	NAG	C4-C5-C6-O6
9	C	1523	NAG	C4-C5-C6-O6
9	C	1542	NAG	C4-C5-C6-O6
9	A	1572	NAG	C4-C5-C6-O6
9	B	1572	NAG	C4-C5-C6-O6
9	C	1573	NAG	C4-C5-C6-O6
9	A	1569	NAG	C4-C5-C6-O6
9	B	1569	NAG	C4-C5-C6-O6
9	C	1570	NAG	C4-C5-C6-O6
9	A	1522	NAG	C8-C7-N2-C2
9	A	1522	NAG	O7-C7-N2-C2
9	A	1523	NAG	C8-C7-N2-C2
9	A	1523	NAG	O7-C7-N2-C2
9	A	1535	NAG	C8-C7-N2-C2
9	A	1541	NAG	C8-C7-N2-C2
9	A	1541	NAG	O7-C7-N2-C2
9	A	1568	NAG	C8-C7-N2-C2
9	A	1568	NAG	O7-C7-N2-C2
9	A	1569	NAG	C8-C7-N2-C2
9	A	1569	NAG	O7-C7-N2-C2
9	A	1572	NAG	C8-C7-N2-C2
9	A	1572	NAG	O7-C7-N2-C2
9	B	1522	NAG	C8-C7-N2-C2
9	B	1522	NAG	O7-C7-N2-C2
9	B	1523	NAG	C8-C7-N2-C2
9	B	1523	NAG	O7-C7-N2-C2
9	B	1535	NAG	C8-C7-N2-C2
9	B	1541	NAG	C8-C7-N2-C2
9	B	1541	NAG	O7-C7-N2-C2
9	B	1568	NAG	C8-C7-N2-C2
9	B	1568	NAG	O7-C7-N2-C2
9	B	1569	NAG	C8-C7-N2-C2
9	B	1569	NAG	O7-C7-N2-C2
9	B	1572	NAG	C8-C7-N2-C2
9	B	1572	NAG	O7-C7-N2-C2
9	C	1523	NAG	C8-C7-N2-C2
9	C	1523	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
9	C	1524	NAG	C8-C7-N2-C2
9	C	1524	NAG	O7-C7-N2-C2
9	C	1536	NAG	C8-C7-N2-C2
9	C	1542	NAG	C8-C7-N2-C2
9	C	1542	NAG	O7-C7-N2-C2
9	C	1569	NAG	C8-C7-N2-C2
9	C	1569	NAG	O7-C7-N2-C2
9	C	1570	NAG	C8-C7-N2-C2
9	C	1570	NAG	O7-C7-N2-C2
9	C	1573	NAG	C8-C7-N2-C2
9	C	1573	NAG	O7-C7-N2-C2
9	A	1568	NAG	C4-C5-C6-O6
9	B	1568	NAG	C4-C5-C6-O6
9	C	1569	NAG	C4-C5-C6-O6
9	A	1572	NAG	O5-C5-C6-O6
9	B	1572	NAG	O5-C5-C6-O6
9	C	1573	NAG	O5-C5-C6-O6
9	A	1535	NAG	O7-C7-N2-C2
9	B	1535	NAG	O7-C7-N2-C2
9	C	1536	NAG	O7-C7-N2-C2
9	A	1568	NAG	O5-C5-C6-O6
9	C	1569	NAG	O5-C5-C6-O6
9	B	1568	NAG	O5-C5-C6-O6
9	A	1522	NAG	C1-C2-N2-C7
9	B	1522	NAG	C1-C2-N2-C7
9	C	1523	NAG	C1-C2-N2-C7
9	A	1522	NAG	C3-C2-N2-C7
9	B	1522	NAG	C3-C2-N2-C7
9	C	1523	NAG	C3-C2-N2-C7

There are no ring outliers.

9 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B	1535	NAG	1	0
9	C	1536	NAG	1	0
9	A	1536	NAG	1	0
9	B	1536	NAG	1	0
10	C	1501	MAN	1	0
9	A	1535	NAG	1	0
10	B	1573	MAN	1	0
9	C	1537	NAG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	A	1573	MAN	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

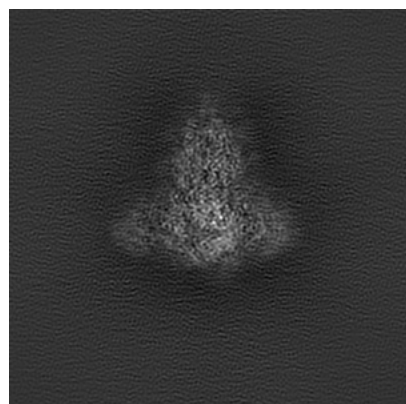
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9891. These allow visual inspection of the internal detail of the map and identification of artifacts.

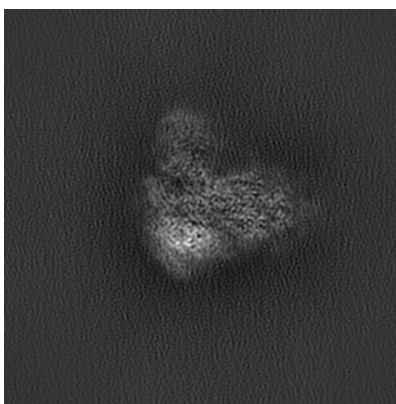
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

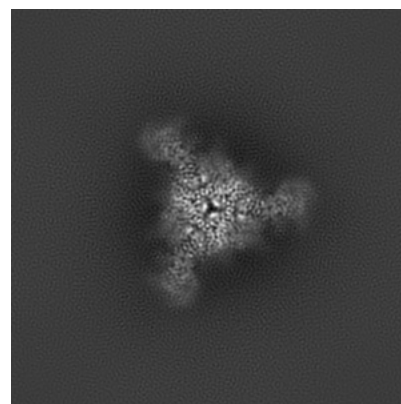
6.1.1 Primary map



X

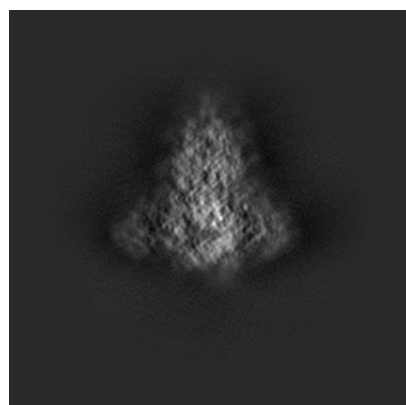


Y

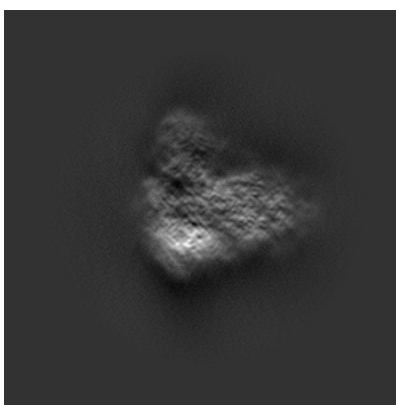


Z

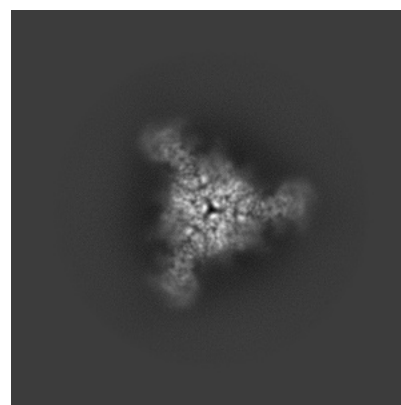
6.1.2 Raw map



X



Y

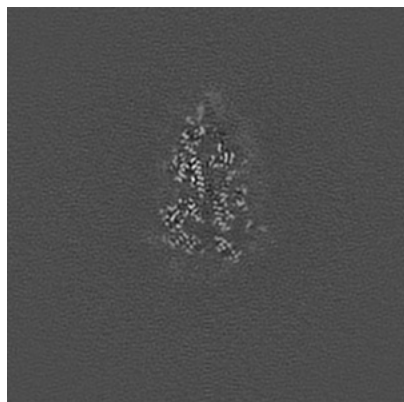


Z

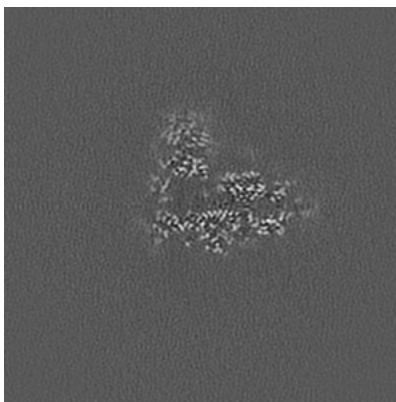
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

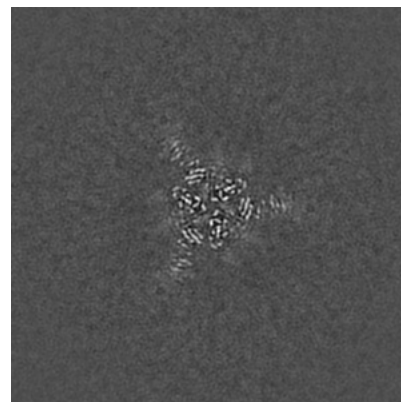
6.2.1 Primary map



X Index: 220

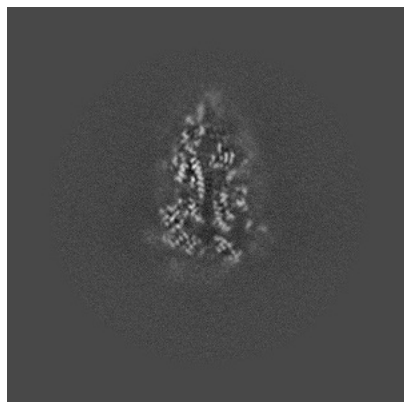


Y Index: 220

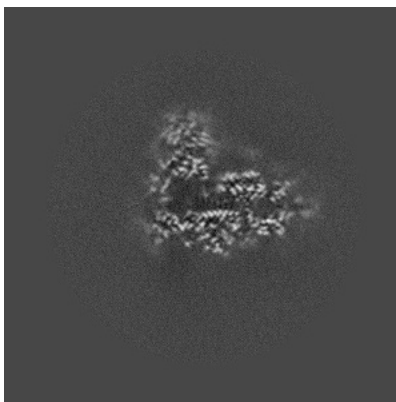


Z Index: 220

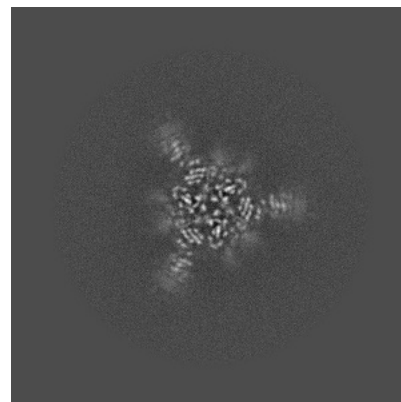
6.2.2 Raw map



X Index: 220



Y Index: 220

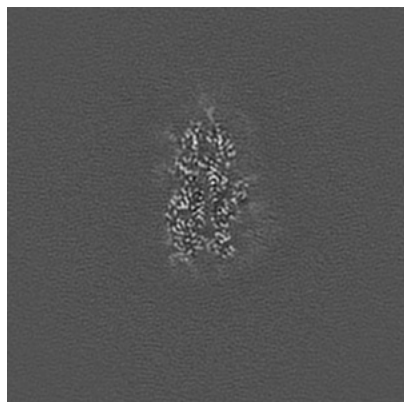


Z Index: 220

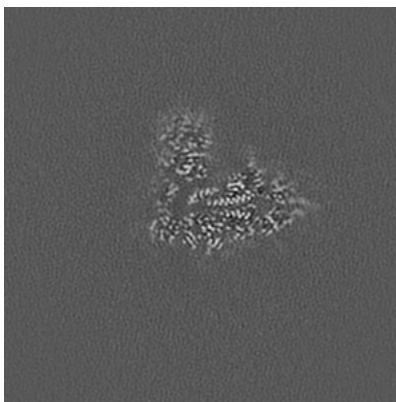
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

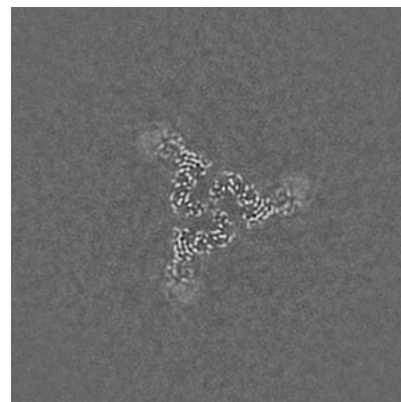
6.3.1 Primary map



X Index: 228

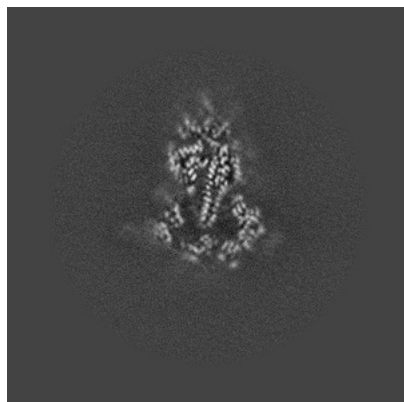


Y Index: 225

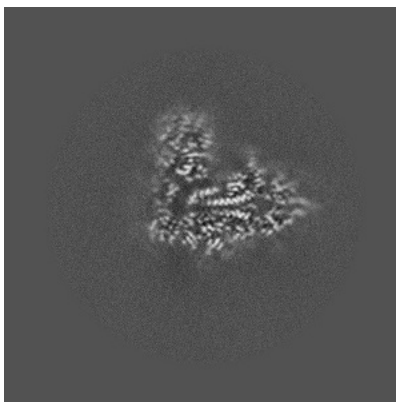


Z Index: 202

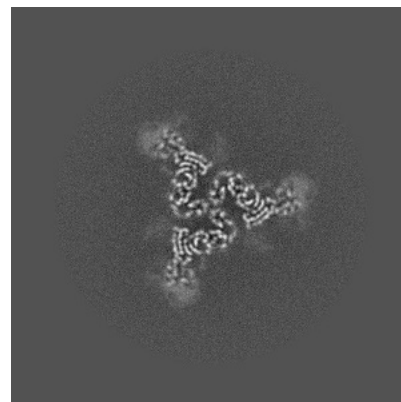
6.3.2 Raw map



X Index: 210



Y Index: 225

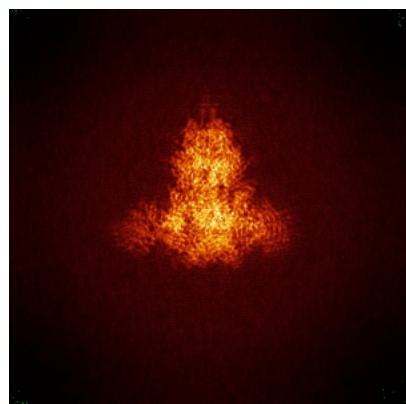


Z Index: 203

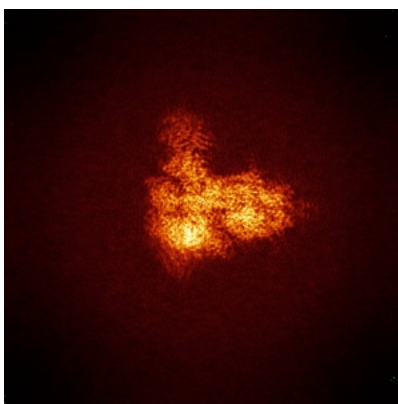
The images above show the largest variance slices of the map in three orthogonal directions.

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

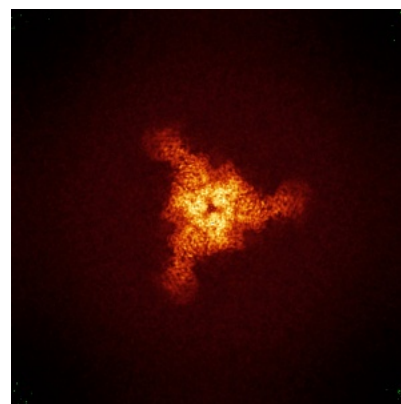
6.4.1 Primary map



X

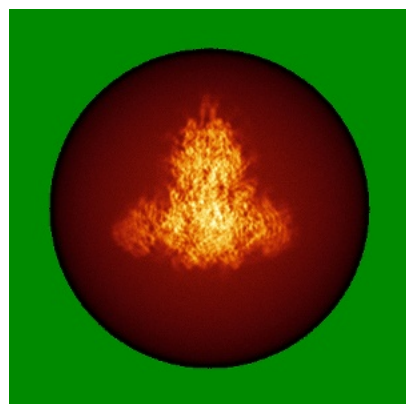


Y

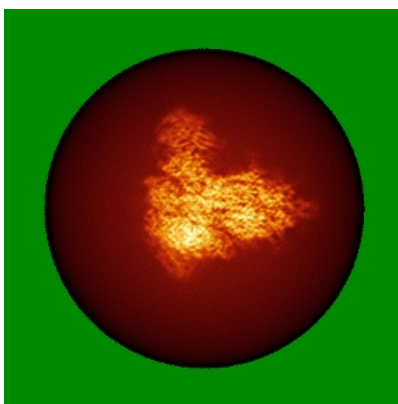


Z

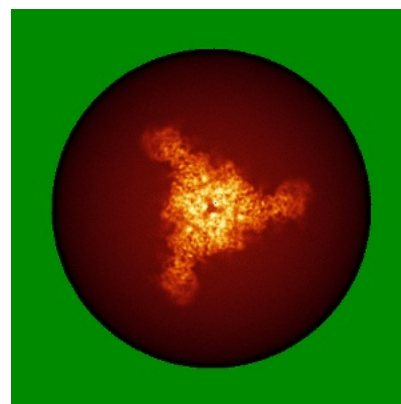
6.4.2 Raw map



X



Y

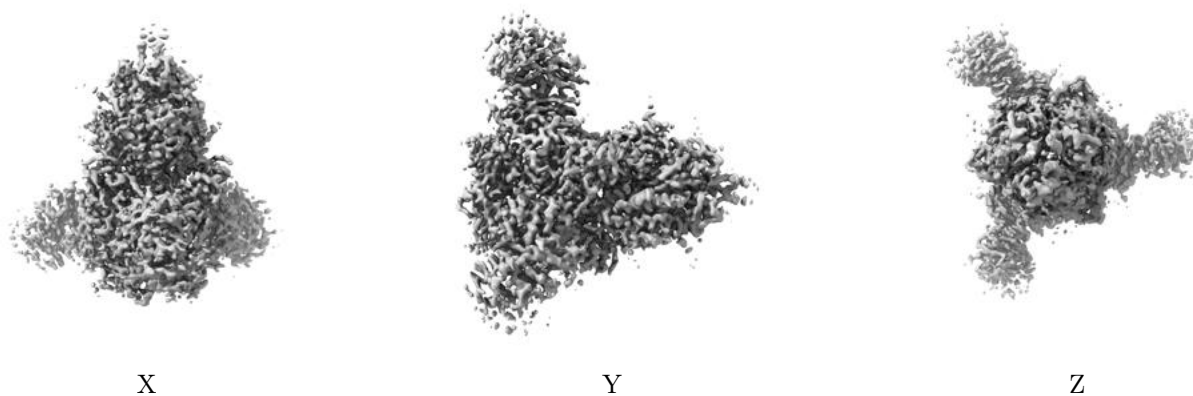


Z

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.

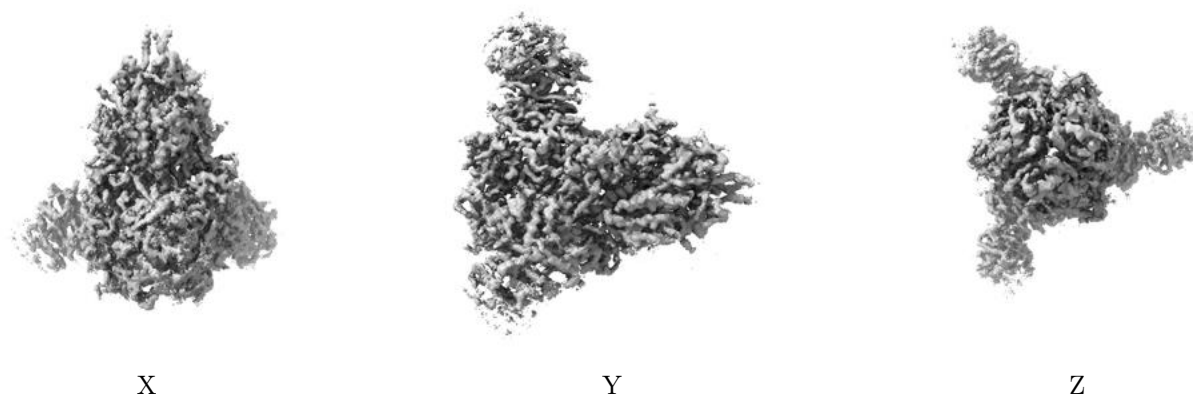
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 5.0. 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.

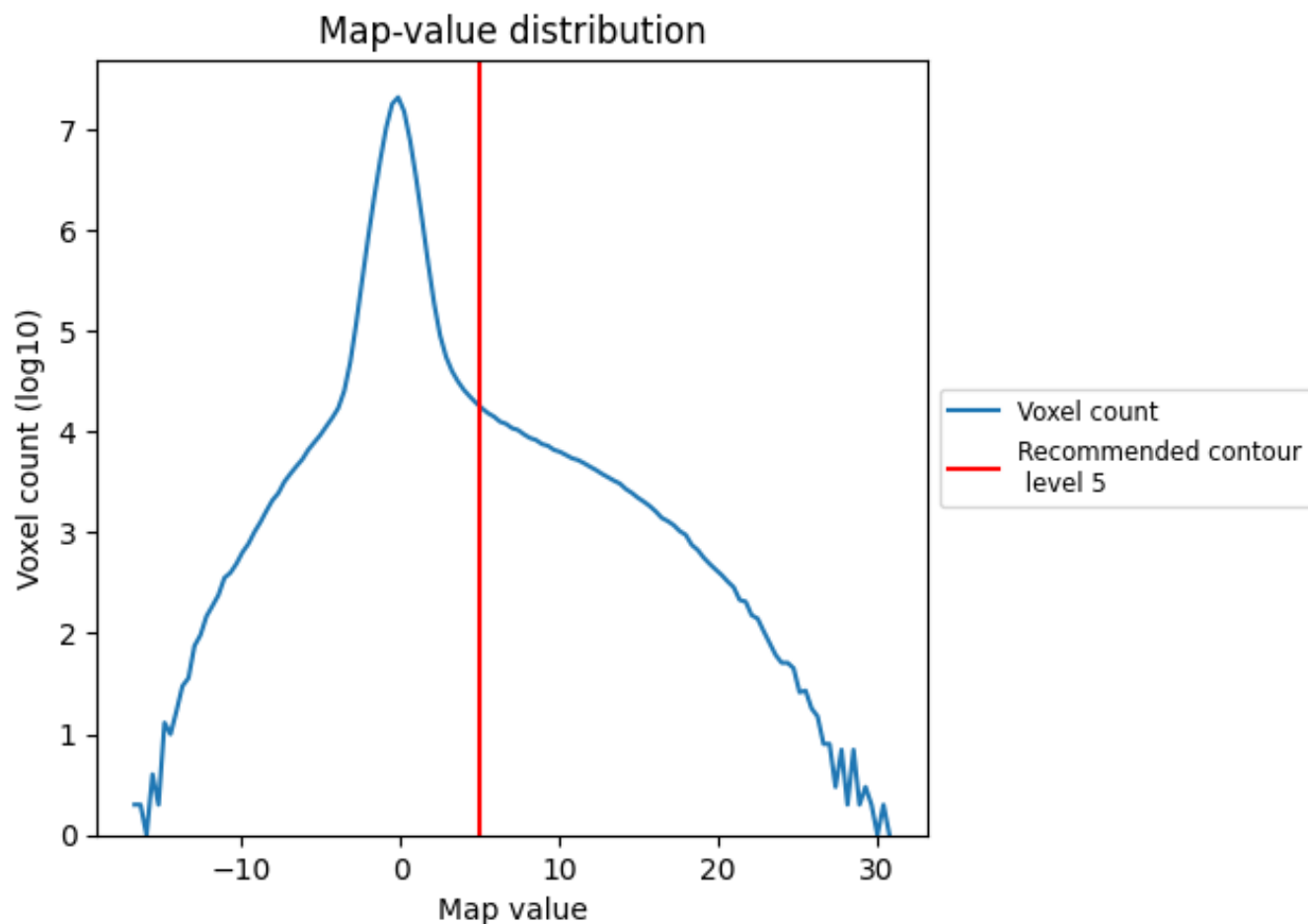
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

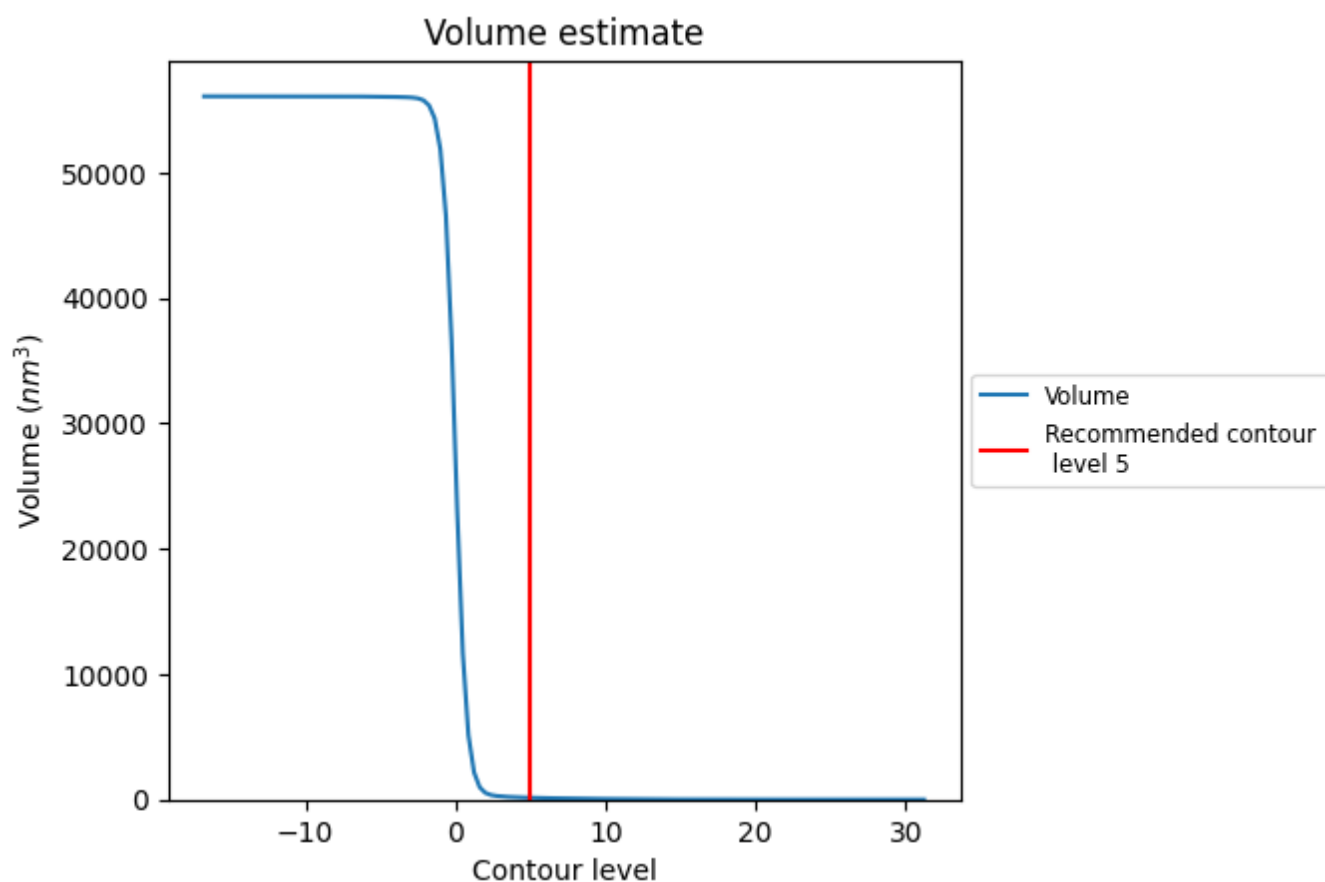
This section contains the results of statistical analysis of the map.

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.

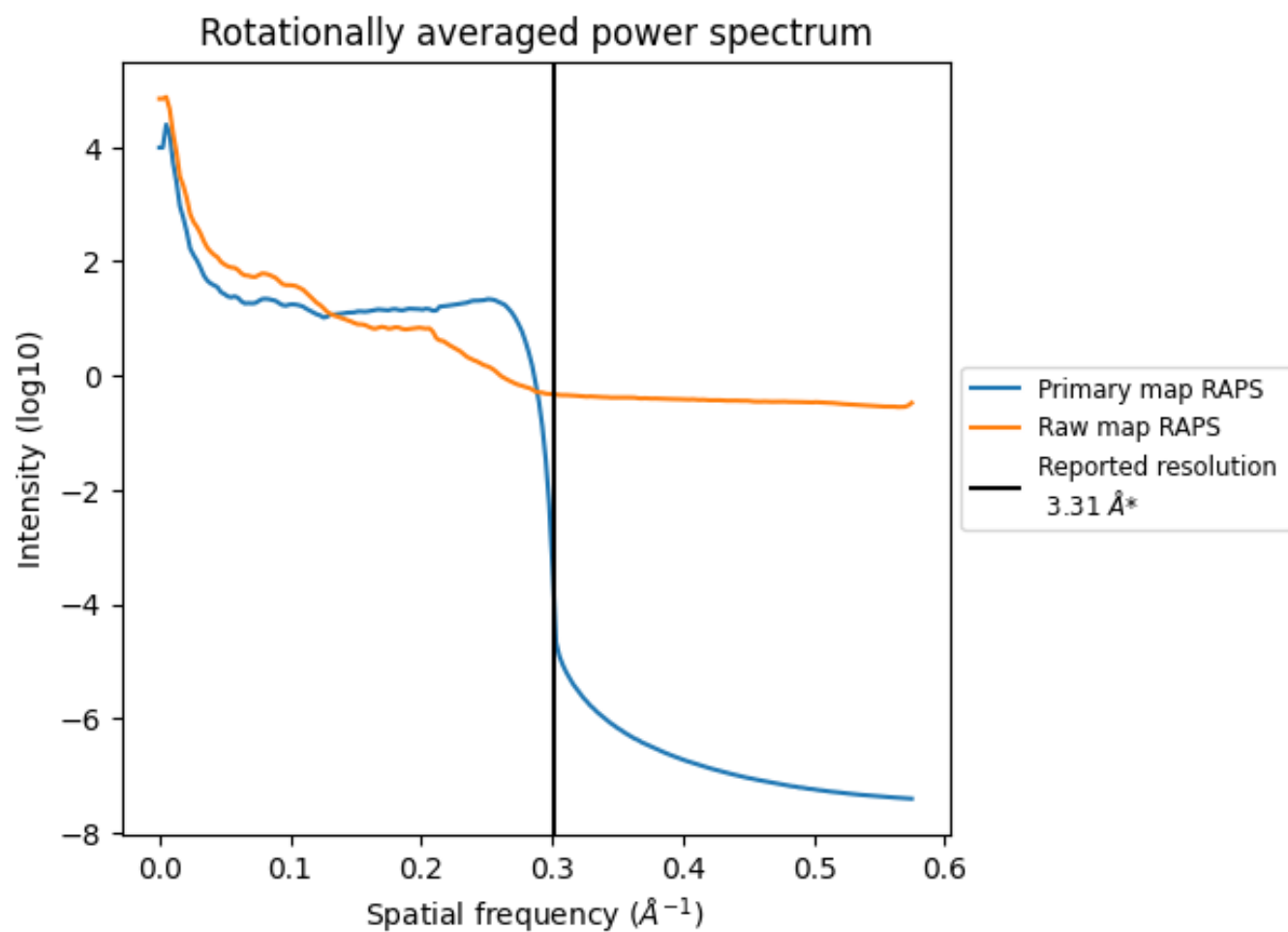
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 145 nm³; this corresponds to an approximate mass of 131 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](#)

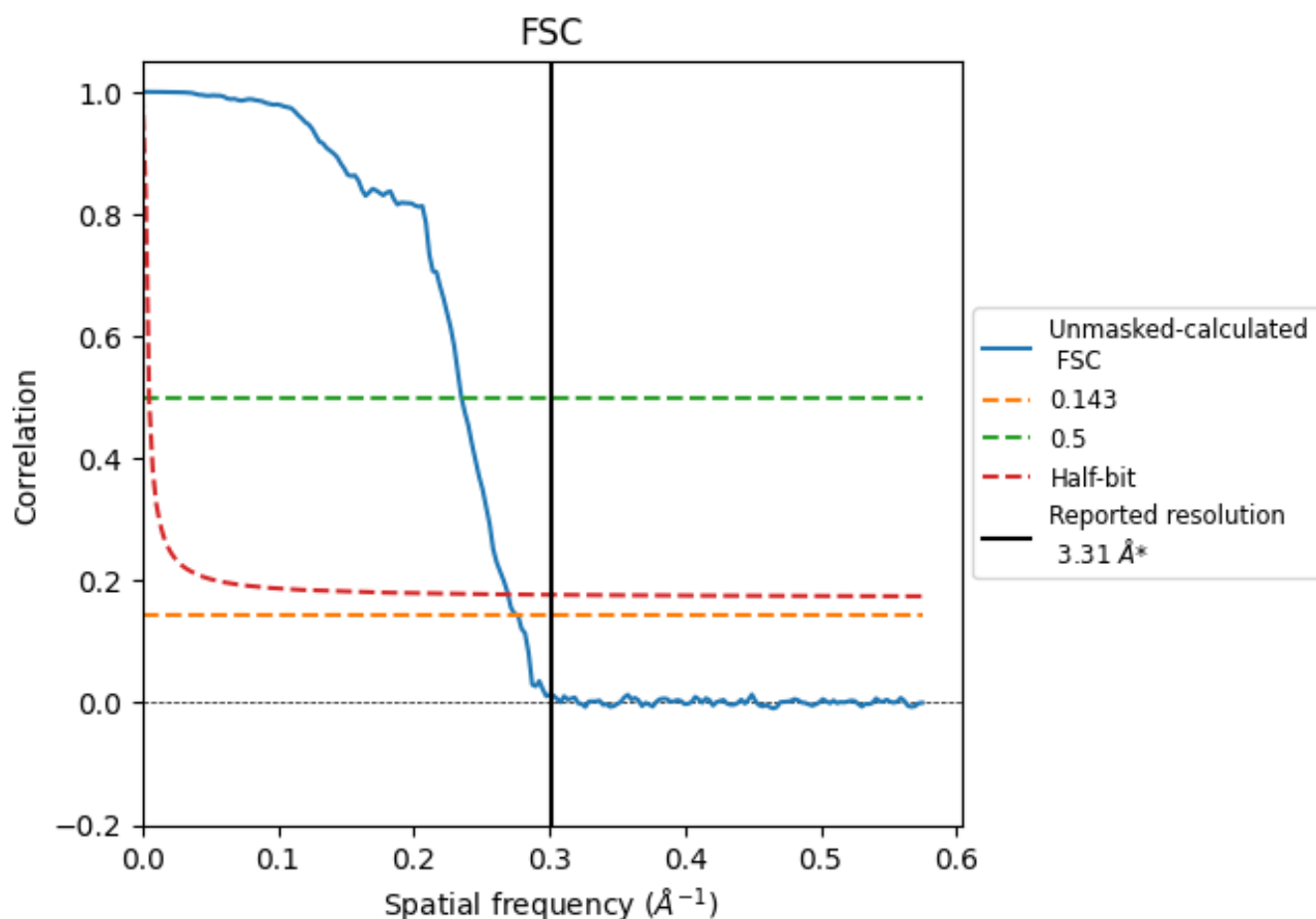


*Reported resolution corresponds to spatial frequency of 0.302 Å⁻¹

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 0.302 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

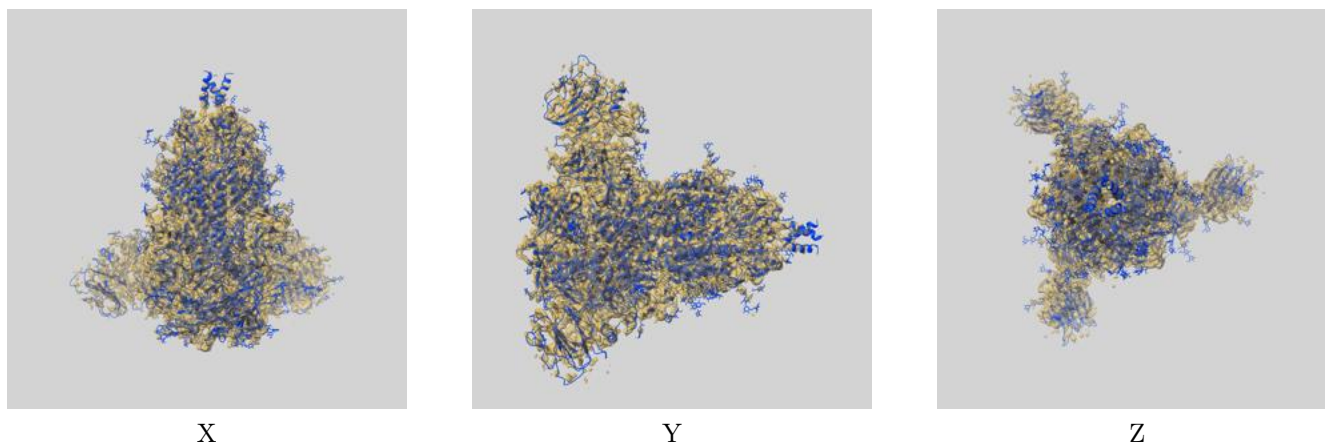
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.31	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.61	4.25	3.71

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

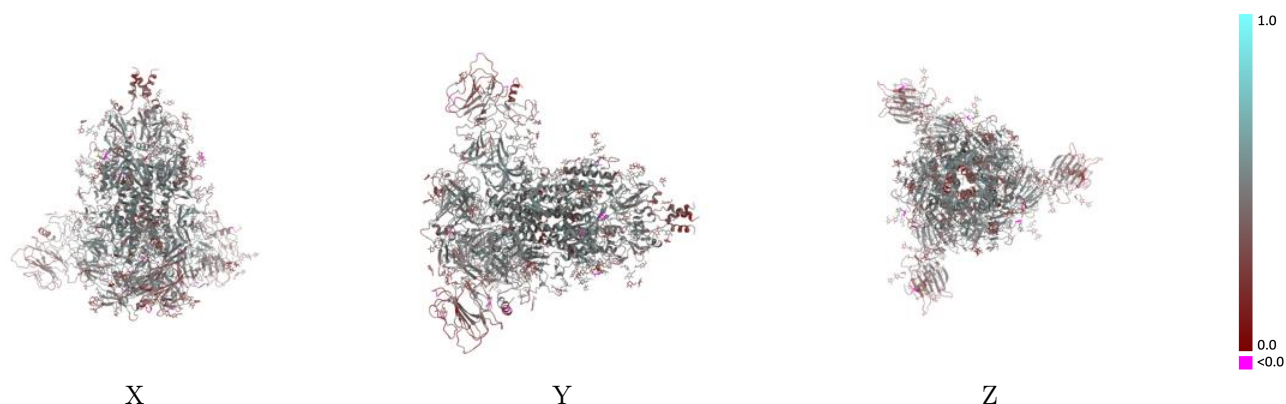
This section contains information regarding the fit between EMDB map EMD-9891 and PDB model 6JX7. Per-residue inclusion information can be found in [section 3](#) on [page 13](#).

9.1 Map-model overlay [i](#)



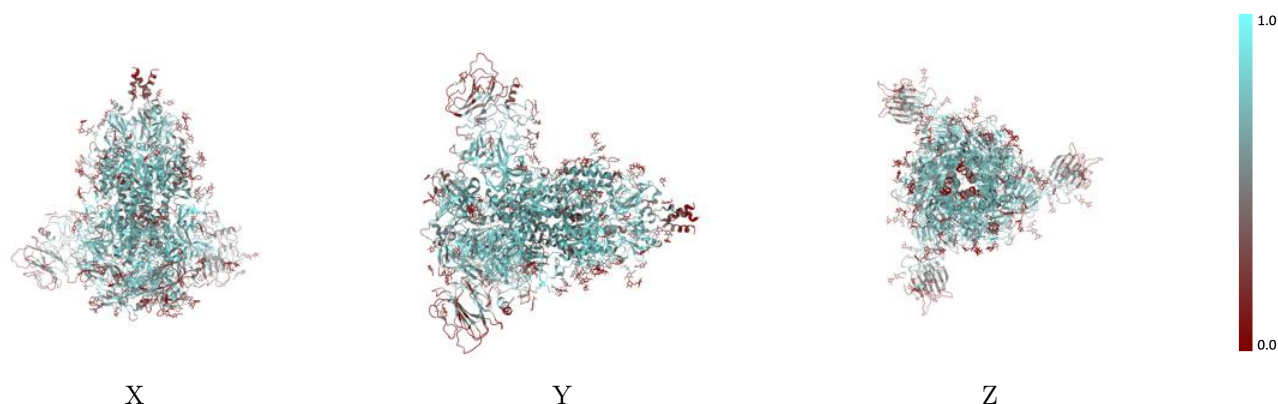
The images above show the 3D surface view of the map at the recommended contour level 5.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



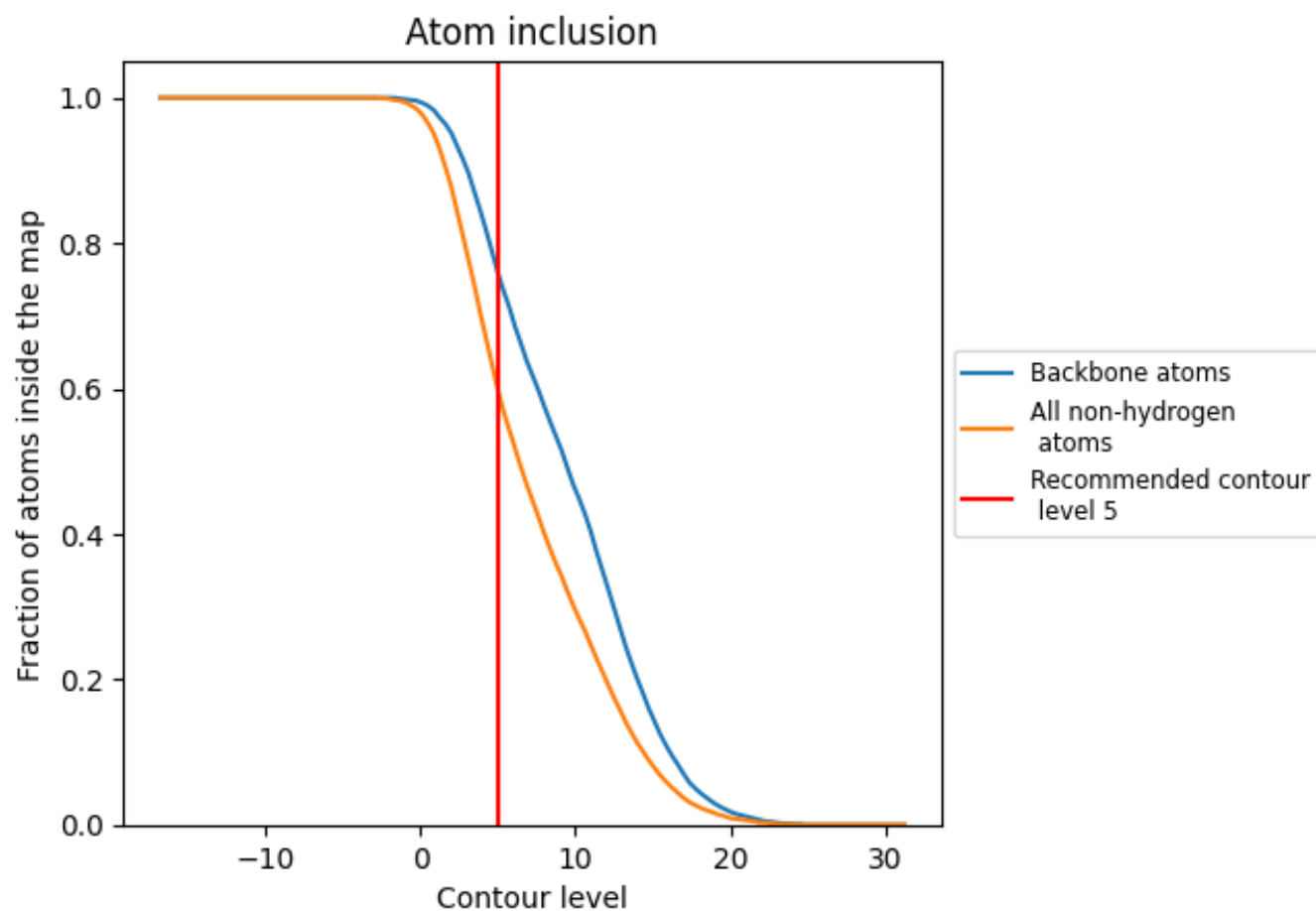
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 60% 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 (5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5980	 0.4340
0	 0.2500	 0.3370
1	 0.2860	 0.3430
2	 0.4100	 0.2910
3	 0.2650	 0.2960
4	 0.2500	 0.3710
A	 0.6250	 0.4410
B	 0.6240	 0.4420
C	 0.6230	 0.4410
D	 0.2500	 0.2860
E	 0.2560	 0.2710
F	 0.3930	 0.2940
G	 0.3080	 0.2780
H	 0.1030	 0.3010
I	 0.3860	 0.3700
J	 0.2400	 0.4400
K	 0.2140	 0.3930
L	 0.2400	 0.2580
M	 0.3570	 0.4580
N	 0.1430	 0.3860
O	 0.2600	 0.4250
P	 0.2400	 0.3420
Q	 0.2500	 0.3360
R	 0.2860	 0.3410
S	 0.3930	 0.2970
T	 0.2770	 0.2920
U	 0.2500	 0.3700
V	 0.2500	 0.3000
W	 0.3080	 0.2680
X	 0.3930	 0.2680
Y	 0.3080	 0.2930
Z	 0.1030	 0.3080
a	 0.3610	 0.3730
b	 0.2400	 0.4400
c	 0.2140	 0.3700



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Chain	Atom inclusion	Q-score
d	 0.2600	 0.2570
e	 0.3570	 0.4640
f	 0.1430	 0.3740
g	 0.2400	 0.4120
h	 0.2400	 0.3460
i	 0.2500	 0.3390
j	 0.2860	 0.3480
k	 0.3930	 0.2960
l	 0.2770	 0.2900
m	 0.2500	 0.3690
n	 0.2500	 0.2810
o	 0.2560	 0.2610
p	 0.3930	 0.2640
q	 0.3080	 0.3010
r	 0.1030	 0.3170
s	 0.3610	 0.3720
t	 0.2400	 0.4410
u	 0.2140	 0.3840
v	 0.2400	 0.2590
w	 0.3210	 0.4470
x	 0.1430	 0.3780
y	 0.2600	 0.4260
z	 0.2600	 0.3420