



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 09:22 PM UTC

PDB ID : 7SBR / pdb_00007sbr
EMDB ID : EMD-24986
Title : One RBD-up 2 of pre-fusion SARS-CoV-2 Kappa variant spike protein
Authors : Zhang, J.; Xiao, T.S.; Cai, Y.F.; Peng, H.Q.; Volloch, S.R.; Chen, B.
Deposited on : 2021-09-25
Resolution : 4.30 Å(reported)
Based on initial model : 7KRR

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:

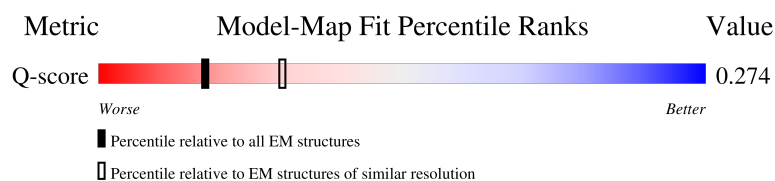
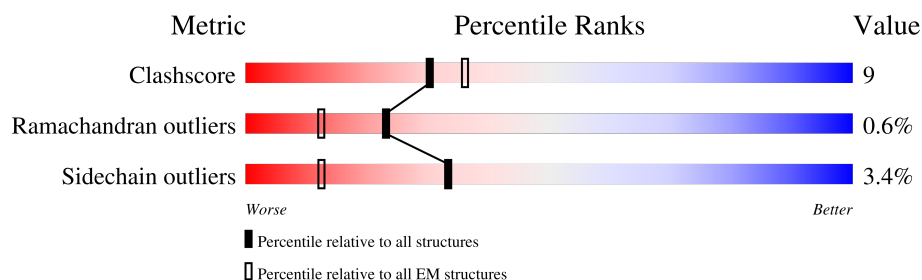
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 4.30 Å.

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



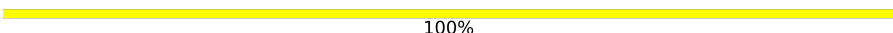
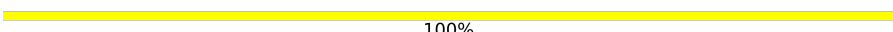
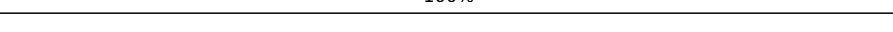
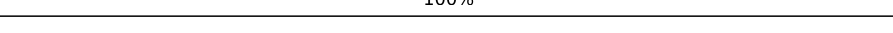
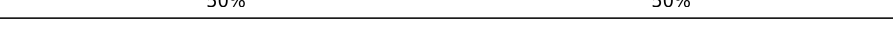
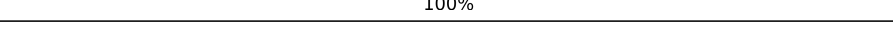
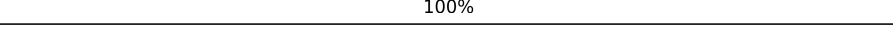
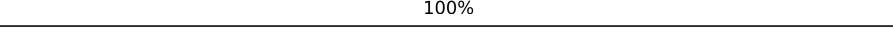
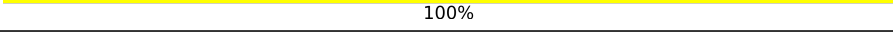
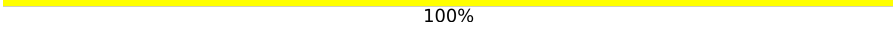
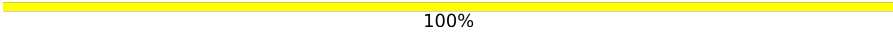
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	4585 (3.80 - 4.80)

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	1310	
1	B	1310	
1	C	1310	
2	D	2	

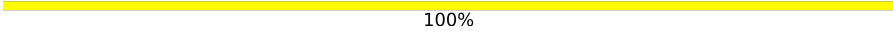
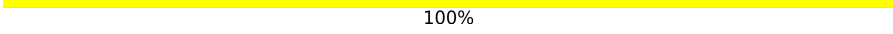
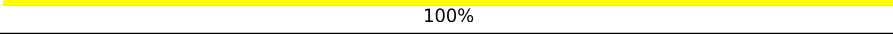
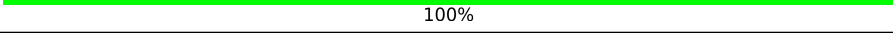
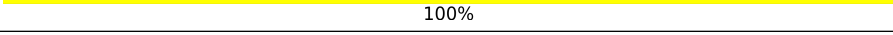
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Mol	Chain	Length	Quality of chain
2	E	2	 50%50%
2	F	2	 100%
2	G	2	 100%
2	H	2	 50%50%
2	N	2	 50%50%
2	O	2	 100%
2	P	2	 50%50%
2	Q	2	 100%
2	R	2	 100%
2	S	2	 100%
2	T	2	 100%
2	Y	2	 50%50%
2	Z	2	 50%50%
2	a	2	 100%
2	b	2	 100%
2	c	2	 50%50%
2	d	2	 100%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

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Mol	Chain	Length	Quality of chain
4	V	3	 67%33%
4	W	3	 33%67%
4	X	3	 100%
4	g	3	 100%
4	h	3	 100%
4	i	3	 100%
4	j	3	 100%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 26959 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 Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1095	Total	C	N	O	S	0	0
			8554	5454	1430	1631	39		
1	B	1076	Total	C	N	O	S	0	0
			8428	5380	1408	1603	37		
1	C	1095	Total	C	N	O	S	0	0
			8566	5465	1433	1630	38		

There are 135 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	154	LYS	GLU	variant	UNP P0DTC2
A	218	HIS	GLN	variant	UNP P0DTC2
A	452	ARG	LEU	variant	UNP P0DTC2
A	484	GLN	GLU	variant	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	681	ARG	PRO	variant	UNP P0DTC2
A	1101	ASP	HIS	variant	UNP P0DTC2
A	1264	LEU	VAL	variant	UNP P0DTC2
A	1274	LEU	-	expression tag	UNP P0DTC2
A	1275	GLU	-	expression tag	UNP P0DTC2
A	1276	SER	-	expression tag	UNP P0DTC2
A	1277	GLY	-	expression tag	UNP P0DTC2
A	1278	GLY	-	expression tag	UNP P0DTC2
A	1279	GLY	-	expression tag	UNP P0DTC2
A	1280	SER	-	expression tag	UNP P0DTC2
A	1281	ALA	-	expression tag	UNP P0DTC2
A	1282	TRP	-	expression tag	UNP P0DTC2
A	1283	SER	-	expression tag	UNP P0DTC2
A	1284	HIS	-	expression tag	UNP P0DTC2
A	1285	PRO	-	expression tag	UNP P0DTC2
A	1286	GLN	-	expression tag	UNP P0DTC2
A	1287	PHE	-	expression tag	UNP P0DTC2
A	1288	GLU	-	expression tag	UNP P0DTC2
A	1289	LYS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1290	GLY	-	expression tag	UNP P0DTC2
A	1291	GLY	-	expression tag	UNP P0DTC2
A	1292	GLY	-	expression tag	UNP P0DTC2
A	1293	SER	-	expression tag	UNP P0DTC2
A	1294	GLY	-	expression tag	UNP P0DTC2
A	1295	GLY	-	expression tag	UNP P0DTC2
A	1296	GLY	-	expression tag	UNP P0DTC2
A	1297	SER	-	expression tag	UNP P0DTC2
A	1298	GLY	-	expression tag	UNP P0DTC2
A	1299	GLY	-	expression tag	UNP P0DTC2
A	1300	SER	-	expression tag	UNP P0DTC2
A	1301	SER	-	expression tag	UNP P0DTC2
A	1302	ALA	-	expression tag	UNP P0DTC2
A	1303	TRP	-	expression tag	UNP P0DTC2
A	1304	SER	-	expression tag	UNP P0DTC2
A	1305	HIS	-	expression tag	UNP P0DTC2
A	1306	PRO	-	expression tag	UNP P0DTC2
A	1307	GLN	-	expression tag	UNP P0DTC2
A	1308	PHE	-	expression tag	UNP P0DTC2
A	1309	GLU	-	expression tag	UNP P0DTC2
A	1310	LYS	-	expression tag	UNP P0DTC2
B	154	LYS	GLU	variant	UNP P0DTC2
B	218	HIS	GLN	variant	UNP P0DTC2
B	452	ARG	LEU	variant	UNP P0DTC2
B	484	GLN	GLU	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	681	ARG	PRO	variant	UNP P0DTC2
B	1101	ASP	HIS	variant	UNP P0DTC2
B	1264	LEU	VAL	variant	UNP P0DTC2
B	1274	LEU	-	expression tag	UNP P0DTC2
B	1275	GLU	-	expression tag	UNP P0DTC2
B	1276	SER	-	expression tag	UNP P0DTC2
B	1277	GLY	-	expression tag	UNP P0DTC2
B	1278	GLY	-	expression tag	UNP P0DTC2
B	1279	GLY	-	expression tag	UNP P0DTC2
B	1280	SER	-	expression tag	UNP P0DTC2
B	1281	ALA	-	expression tag	UNP P0DTC2
B	1282	TRP	-	expression tag	UNP P0DTC2
B	1283	SER	-	expression tag	UNP P0DTC2
B	1284	HIS	-	expression tag	UNP P0DTC2
B	1285	PRO	-	expression tag	UNP P0DTC2
B	1286	GLN	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1287	PHE	-	expression tag	UNP P0DTC2
B	1288	GLU	-	expression tag	UNP P0DTC2
B	1289	LYS	-	expression tag	UNP P0DTC2
B	1290	GLY	-	expression tag	UNP P0DTC2
B	1291	GLY	-	expression tag	UNP P0DTC2
B	1292	GLY	-	expression tag	UNP P0DTC2
B	1293	SER	-	expression tag	UNP P0DTC2
B	1294	GLY	-	expression tag	UNP P0DTC2
B	1295	GLY	-	expression tag	UNP P0DTC2
B	1296	GLY	-	expression tag	UNP P0DTC2
B	1297	SER	-	expression tag	UNP P0DTC2
B	1298	GLY	-	expression tag	UNP P0DTC2
B	1299	GLY	-	expression tag	UNP P0DTC2
B	1300	SER	-	expression tag	UNP P0DTC2
B	1301	SER	-	expression tag	UNP P0DTC2
B	1302	ALA	-	expression tag	UNP P0DTC2
B	1303	TRP	-	expression tag	UNP P0DTC2
B	1304	SER	-	expression tag	UNP P0DTC2
B	1305	HIS	-	expression tag	UNP P0DTC2
B	1306	PRO	-	expression tag	UNP P0DTC2
B	1307	GLN	-	expression tag	UNP P0DTC2
B	1308	PHE	-	expression tag	UNP P0DTC2
B	1309	GLU	-	expression tag	UNP P0DTC2
B	1310	LYS	-	expression tag	UNP P0DTC2
C	154	LYS	GLU	variant	UNP P0DTC2
C	218	HIS	GLN	variant	UNP P0DTC2
C	452	ARG	LEU	variant	UNP P0DTC2
C	484	GLN	GLU	variant	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	681	ARG	PRO	variant	UNP P0DTC2
C	1101	ASP	HIS	variant	UNP P0DTC2
C	1264	LEU	VAL	variant	UNP P0DTC2
C	1274	LEU	-	expression tag	UNP P0DTC2
C	1275	GLU	-	expression tag	UNP P0DTC2
C	1276	SER	-	expression tag	UNP P0DTC2
C	1277	GLY	-	expression tag	UNP P0DTC2
C	1278	GLY	-	expression tag	UNP P0DTC2
C	1279	GLY	-	expression tag	UNP P0DTC2
C	1280	SER	-	expression tag	UNP P0DTC2
C	1281	ALA	-	expression tag	UNP P0DTC2
C	1282	TRP	-	expression tag	UNP P0DTC2
C	1283	SER	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1284	HIS	-	expression tag	UNP P0DTC2
C	1285	PRO	-	expression tag	UNP P0DTC2
C	1286	GLN	-	expression tag	UNP P0DTC2
C	1287	PHE	-	expression tag	UNP P0DTC2
C	1288	GLU	-	expression tag	UNP P0DTC2
C	1289	LYS	-	expression tag	UNP P0DTC2
C	1290	GLY	-	expression tag	UNP P0DTC2
C	1291	GLY	-	expression tag	UNP P0DTC2
C	1292	GLY	-	expression tag	UNP P0DTC2
C	1293	SER	-	expression tag	UNP P0DTC2
C	1294	GLY	-	expression tag	UNP P0DTC2
C	1295	GLY	-	expression tag	UNP P0DTC2
C	1296	GLY	-	expression tag	UNP P0DTC2
C	1297	SER	-	expression tag	UNP P0DTC2
C	1298	GLY	-	expression tag	UNP P0DTC2
C	1299	GLY	-	expression tag	UNP P0DTC2
C	1300	SER	-	expression tag	UNP P0DTC2
C	1301	SER	-	expression tag	UNP P0DTC2
C	1302	ALA	-	expression tag	UNP P0DTC2
C	1303	TRP	-	expression tag	UNP P0DTC2
C	1304	SER	-	expression tag	UNP P0DTC2
C	1305	HIS	-	expression tag	UNP P0DTC2
C	1306	PRO	-	expression tag	UNP P0DTC2
C	1307	GLN	-	expression tag	UNP P0DTC2
C	1308	PHE	-	expression tag	UNP P0DTC2
C	1309	GLU	-	expression tag	UNP P0DTC2
C	1310	LYS	-	expression tag	UNP P0DTC2

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



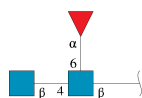
Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	2	Total	C	N	O	0	0
			28	16	2	10		
2	E	2	Total	C	N	O	0	0
			28	16	2	10		
2	F	2	Total	C	N	O	0	0
			28	16	2	10		

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Mol	Chain	Residues	Atoms				AltConf	Trace
2	G	2	Total	C	N	O	0	0
			28	16	2	10		
2	H	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	O	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	Q	2	Total	C	N	O	0	0
			28	16	2	10		
2	R	2	Total	C	N	O	0	0
			28	16	2	10		
2	S	2	Total	C	N	O	0	0
			28	16	2	10		
2	T	2	Total	C	N	O	0	0
			28	16	2	10		
2	Y	2	Total	C	N	O	0	0
			28	16	2	10		
2	Z	2	Total	C	N	O	0	0
			28	16	2	10		
2	a	2	Total	C	N	O	0	0
			28	16	2	10		
2	b	2	Total	C	N	O	0	0
			28	16	2	10		
2	c	2	Total	C	N	O	0	0
			28	16	2	10		
2	d	2	Total	C	N	O	0	0
			28	16	2	10		
2	e	2	Total	C	N	O	0	0
			28	16	2	10		

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



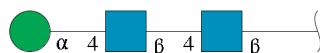
Mol	Chain	Residues	Atoms				AltConf	Trace
3	I	3	Total	C	N	O	0	0
			38	22	2	14		

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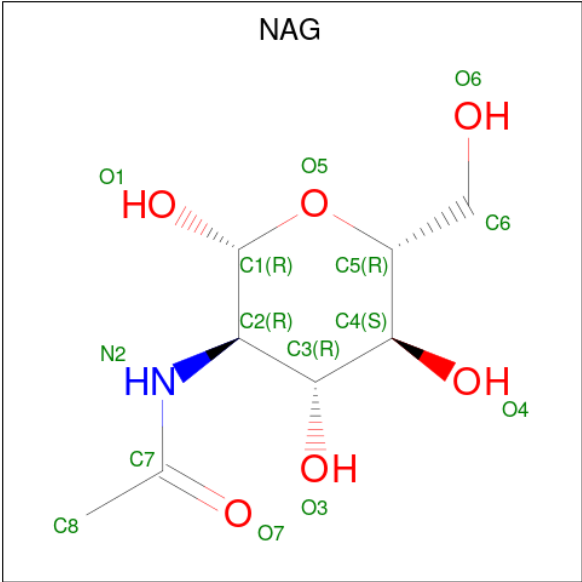
Mol	Chain	Residues	Atoms				AltConf	Trace
3	U	3	Total	C	N	O	0	0
			38	22	2	14		
3	f	3	Total	C	N	O	0	0
			38	22	2	14		

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



Mol	Chain	Residues	Atoms				AltConf	Trace
4	J	3	Total	C	N	O	0	0
			39	22	2	15		
4	K	3	Total	C	N	O	0	0
			39	22	2	15		
4	L	3	Total	C	N	O	0	0
			39	22	2	15		
4	M	3	Total	C	N	O	0	0
			39	22	2	15		
4	V	3	Total	C	N	O	0	0
			39	22	2	15		
4	W	3	Total	C	N	O	0	0
			39	22	2	15		
4	X	3	Total	C	N	O	0	0
			39	22	2	15		
4	g	3	Total	C	N	O	0	0
			39	22	2	15		
4	h	3	Total	C	N	O	0	0
			39	22	2	15		
4	i	3	Total	C	N	O	0	0
			39	22	2	15		
4	j	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).

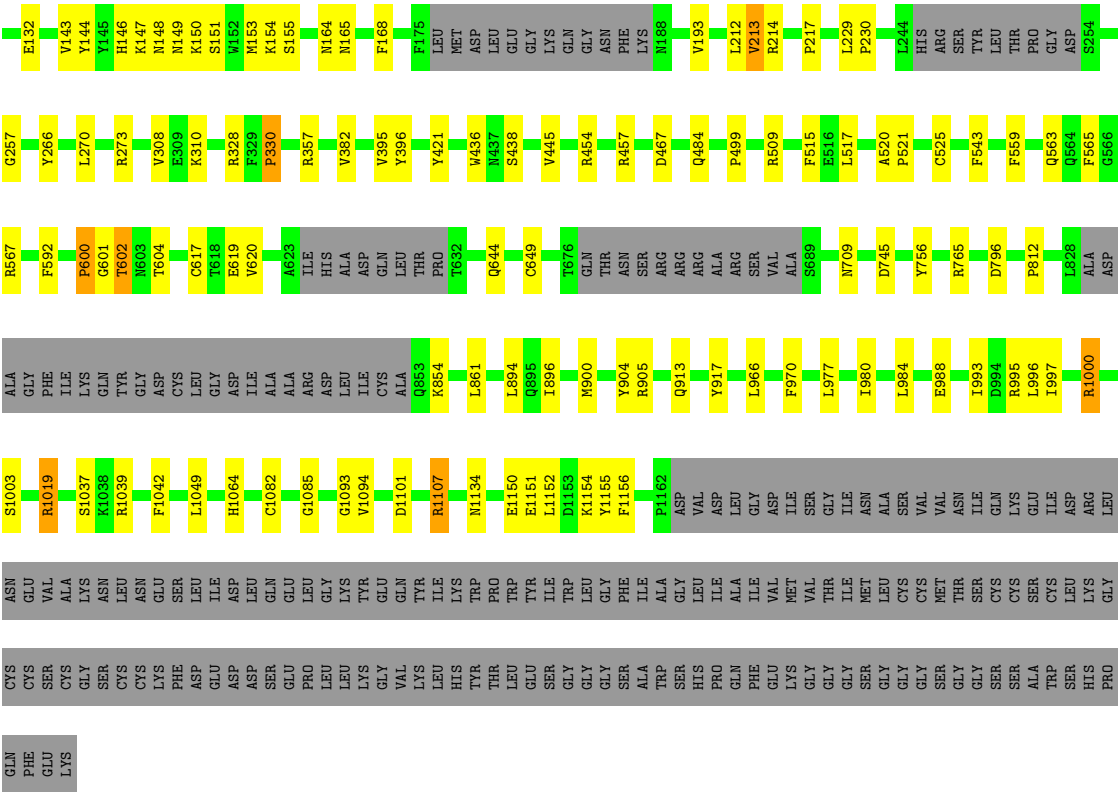


Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5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

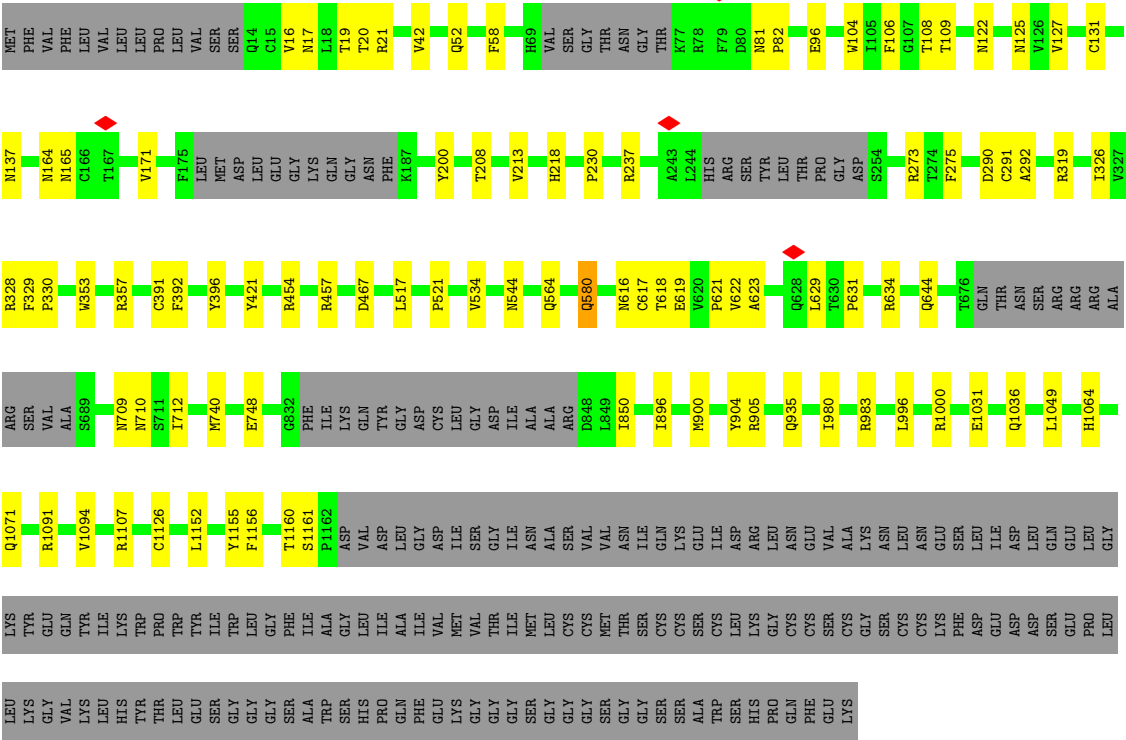
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Mol	Chain	Residues	Atoms				AltConf
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C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



• Molecule 1: Spike glycoprotein



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

Chain D:  50% 50%

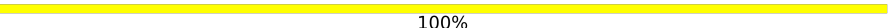
MAG1
MAG2

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

Chain E:  50% 50%

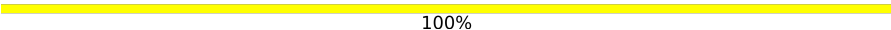
MAG1
MAG2

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

Chain F:  100%

MAG1
MAG2

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

Chain G:  100%

MAG1
MAG2

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

Chain H:  50% 50%

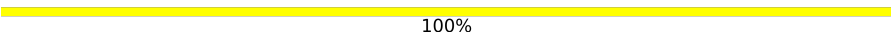
MAG1
MAG2

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

Chain N:  50% 50%

MAG1
MAG2

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

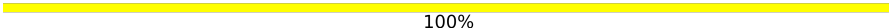
Chain O:  100%

MAG1
MAG2

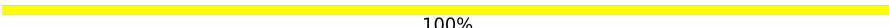
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:  50% 50%MAG1
MAG2

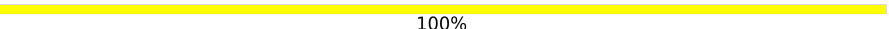
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  100%MAG1
MAG2

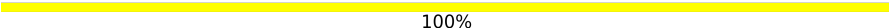
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain R:  100%MAG1
MAG2

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

Chain S:  100%MAG1
MAG2

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

Chain T:  100%MAG1
MAG2

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

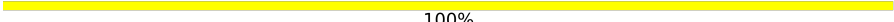
Chain Y:  50% 50% 50%MAG1
MAG2

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

Chain Z:  50% 50%

MAG1
MAG2

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

Chain a:  100%

MAG1
MAG2

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

Chain b:  100%

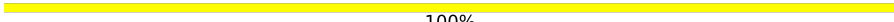
MAG1
MAG2

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

Chain c:  50% 50%

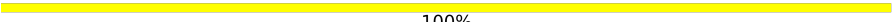
MAG1
MAG2

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

Chain d:  100%

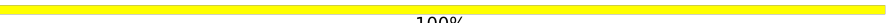
MAG1
MAG2

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

Chain e:  100%

MAG1
MAG2

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

Chain I:  100%

MAG1
MAG2
FUC3

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

Chain U:  33% 67%

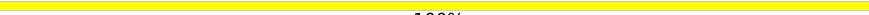
MAG1
MAG2
FUC3

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

Chain f:  100%

MAG1
MAG2
FUC3

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

Chain J:  100%

MAG1
MAG2
MAN3

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

Chain K:  33% 67%

MAG1
MAG2
MAN3

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

Chain L:  100%

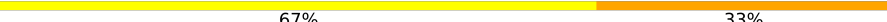
MAG1
MAG2
MAN3

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

Chain M:  100%

MAG1
MAG2
MAN3

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

Chain V:  67% 33%

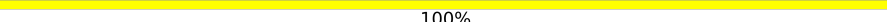
MAG1
MAG2
MAN3

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

Chain W:  33% 67%

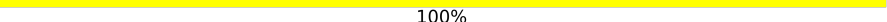
MAG1
MAG2
MAN3

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

Chain X:  100%

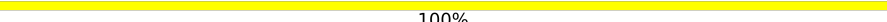
MAG1
MAG2
MAN3

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

Chain g:  100%

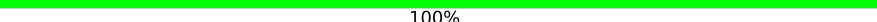
MAG1
MAG2
MAN3

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

Chain h:  100%

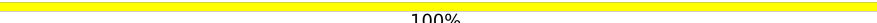
MAG1
MAG2
MAN3

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

Chain i:  100%

MAG1
MAG2
MAN3

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

Chain j:  100%

MAG1
MAG2
MAG3

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	21830	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	51.63	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.332	Depositor
Minimum map value	-0.002	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.052	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	396.03217, 396.03217, 396.03217	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.825067, 0.825067, 0.825067	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.74	3/8749 (0.0%)	1.28	14/11900 (0.1%)
1	B	0.71	1/8623 (0.0%)	1.28	22/11731 (0.2%)
1	C	0.72	0/8765	1.28	16/11927 (0.1%)
All	All	0.72	4/26137 (0.0%)	1.28	52/35558 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1134	ASN	C-N	-6.80	1.24	1.33
1	A	44	ARG	CA-C	6.15	1.59	1.52
1	A	1000	ARG	N-CA	-5.80	1.39	1.46
1	B	1000	ARG	N-CA	-5.55	1.39	1.46

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	164	ASN	N-CA-C	10.23	123.38	111.11
1	A	1107	ARG	N-CA-C	8.85	123.84	112.34
1	B	213	VAL	N-CA-C	8.82	124.38	112.04
1	B	1107	ARG	N-CA-C	8.20	123.00	112.34
1	C	1091	ARG	N-CA-C	-7.84	102.67	111.14
1	A	102	ARG	N-CA-C	7.38	122.46	113.38
1	B	122	ASN	CA-CB-CG	7.32	119.92	112.60
1	A	565	PHE	CA-CB-CG	7.13	120.93	113.80
1	B	602	THR	N-CA-C	-6.82	103.84	111.28
1	B	484	GLN	CA-C-N	6.75	130.94	121.09
1	B	484	GLN	C-N-CA	6.75	130.94	121.09
1	A	830	ASP	CA-CB-CG	6.53	119.13	112.60
1	A	747	THR	CA-C-N	6.51	128.90	120.44
1	A	747	THR	C-N-CA	6.51	128.90	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	559	PHE	CA-CB-CG	6.25	120.05	113.80
1	A	559	PHE	CA-CB-CG	6.09	119.89	113.80
1	C	1161	SER	N-CA-C	6.01	120.40	112.35
1	B	330	PRO	N-CA-CB	5.99	109.02	103.39
1	B	96	GLU	N-CA-C	5.86	118.31	109.23
1	B	565	PHE	CA-C-N	5.85	126.11	121.61
1	B	565	PHE	C-N-CA	5.85	126.11	121.61
1	A	481	ASN	CA-CB-CG	5.82	118.42	112.60
1	C	16	VAL	N-CA-C	5.80	116.00	108.35
1	C	208	THR	N-CA-C	5.75	117.93	110.40
1	B	214	ARG	N-CA-C	5.74	123.02	110.80
1	C	580	GLN	OE1-CD-NE2	-5.69	116.91	122.60
1	B	745	ASP	CA-CB-CG	5.68	118.28	112.60
1	B	122	ASN	CB-CA-C	5.67	119.76	109.83
1	C	1071	GLN	OE1-CD-NE2	-5.58	117.02	122.60
1	C	850	ILE	CA-C-N	5.57	128.20	120.29
1	C	850	ILE	C-N-CA	5.57	128.20	120.29
1	B	17	ASN	CB-CA-C	-5.50	101.73	110.81
1	A	318	PHE	CA-CB-CG	5.47	119.27	113.80
1	B	1064	HIS	CB-CG-CD2	-5.43	124.14	131.20
1	C	213	VAL	N-CA-C	5.42	115.55	110.30
1	C	1064	HIS	CB-CG-CD2	-5.41	124.16	131.20
1	C	617	CYS	N-CA-C	5.41	118.67	111.75
1	A	1135	ASN	CA-CB-CG	-5.34	107.26	112.60
1	B	16	VAL	O-C-N	5.28	128.91	123.05
1	B	1085	GLY	CA-C-N	5.26	128.69	120.75
1	B	1085	GLY	C-N-CA	5.26	128.69	120.75
1	B	709	ASN	CA-CB-CG	5.21	117.81	112.60
1	A	804	GLN	OE1-CD-NE2	-5.19	117.41	122.60
1	B	1019	ARG	N-CA-C	-5.17	105.55	111.14
1	C	96	GLU	N-CA-CB	-5.17	103.03	111.66
1	C	935	GLN	OE1-CD-NE2	-5.14	117.46	122.60
1	A	1088	HIS	CB-CG-CD2	-5.09	124.58	131.20
1	C	631	PRO	N-CA-CB	5.08	107.36	103.30
1	C	218	HIS	CB-CG-CD2	-5.04	124.65	131.20
1	B	563	GLN	OE1-CD-NE2	-5.03	117.57	122.60
1	A	564	GLN	OE1-CD-NE2	-5.03	117.58	122.60
1	A	944	ALA	N-CA-C	5.02	116.75	111.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	8554	0	8319	214	0
1	B	8428	0	8202	223	0
1	C	8566	0	8335	150	0
2	D	28	0	25	1	0
2	E	28	0	25	1	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	H	28	0	25	1	0
2	N	28	0	25	4	0
2	O	28	0	25	0	0
2	P	28	0	25	1	0
2	Q	28	0	25	0	0
2	R	28	0	25	1	0
2	S	28	0	25	0	0
2	T	28	0	25	0	0
2	Y	28	0	25	1	0
2	Z	28	0	25	4	0
2	a	28	0	25	0	0
2	b	28	0	25	1	0
2	c	28	0	25	2	0
2	d	28	0	25	0	0
2	e	28	0	25	0	0
3	I	38	0	34	0	0
3	U	38	0	34	0	0
3	f	38	0	34	0	0
4	J	39	0	34	0	0
4	K	39	0	34	0	0
4	L	39	0	34	0	0
4	M	39	0	34	0	0
4	V	39	0	34	1	0
4	W	39	0	34	0	0
4	X	39	0	34	0	0
4	g	39	0	34	0	0
4	h	39	0	34	0	0
4	i	39	0	34	0	0
4	j	39	0	34	0	0
5	A	126	0	117	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	112	0	104	6	0
5	C	98	0	91	2	0
All	All	26959	0	26119	475	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 (475) 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:C:58:PHE:CE2	1:C:275:PHE:HE2	1.41	1.37
1:C:58:PHE:CE2	1:C:275:PHE:CE2	2.20	1.29
1:C:58:PHE:CD2	1:C:290:ASP:HB2	1.66	1.27
1:A:437:ASN:HA	1:A:508:TYR:CD1	1.78	1.19
1:A:1079:PRO:HB2	1:B:917:TYR:CE1	1.78	1.18
1:A:124:THR:HB	5:A:1402:NAG:C8	1.76	1.15
1:A:437:ASN:HB2	1:A:508:TYR:HE1	1.03	1.15
1:A:567:ARG:HD2	1:B:42:VAL:HG11	1.28	1.14
1:B:601:GLY:HA3	1:B:604:THR:HG22	1.25	1.12
1:C:58:PHE:HE2	1:C:275:PHE:CE2	1.57	1.11
1:A:1156:PHE:CZ	1:B:1155:TYR:HB3	1.85	1.10
1:C:329:PHE:CE1	1:C:544:ASN:HA	1.87	1.10
1:A:815:ARG:NH2	1:A:867:ASP:HB2	1.67	1.10
1:A:437:ASN:HB2	1:A:508:TYR:CE1	1.86	1.08
1:A:1079:PRO:CB	1:B:917:TYR:HE1	1.66	1.07
1:C:273:ARG:HH12	1:C:292:ALA:HB3	1.09	1.07
1:A:124:THR:CB	5:A:1402:NAG:H82	1.82	1.07
1:C:171:VAL:HG21	2:Z:2:NAG:H82	1.38	1.06
1:A:437:ASN:CA	1:A:508:TYR:CD1	2.39	1.05
1:A:437:ASN:CB	1:A:508:TYR:CE1	2.40	1.04
1:B:57:PRO:HB3	1:B:273:ARG:HH22	1.20	1.04
1:B:601:GLY:HA3	1:B:604:THR:CG2	1.89	1.03
1:C:900:MET:HG3	1:C:904:TYR:HE2	1.22	1.02
1:B:900:MET:HG3	1:B:904:TYR:HE2	1.19	1.02
1:B:900:MET:HG3	1:B:904:TYR:CE2	1.95	1.01
1:A:815:ARG:NH2	1:A:867:ASP:CB	2.22	1.01
1:C:58:PHE:HD2	1:C:290:ASP:HB2	1.05	1.00
1:C:905:ARG:HH11	1:C:1036:GLN:HB2	1.20	1.00
1:A:815:ARG:HH21	1:A:867:ASP:HB2	1.25	1.00
1:A:1079:PRO:CB	1:B:917:TYR:CE1	2.41	1.00
1:A:897:PRO:HD2	1:A:900:MET:HE2	1.42	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:905:ARG:HH11	1:A:1036:GLN:HB2	1.24	0.99
1:C:329:PHE:CE1	1:C:544:ASN:CA	2.46	0.99
1:A:904:TYR:CD1	1:C:1107:ARG:HD3	1.97	0.98
1:C:273:ARG:NH1	1:C:292:ALA:HB3	1.77	0.98
1:A:1107:ARG:HD3	1:B:904:TYR:CD1	1.96	0.98
1:C:980:ILE:HD12	1:C:996:LEU:HD11	1.39	0.98
1:B:132:GLU:HG3	1:B:165:ASN:HB2	1.45	0.97
1:A:796:ASP:OD2	1:C:709:ASN:HB3	1.64	0.97
1:B:395:VAL:HG22	1:B:515:PHE:HB3	1.45	0.96
1:B:567:ARG:NH1	1:C:42:VAL:HG21	1.81	0.95
1:A:966:LEU:HD23	1:A:1000:ARG:HD3	1.49	0.95
1:C:905:ARG:NH1	1:C:1036:GLN:HB2	1.82	0.94
1:A:1107:ARG:CZ	1:B:904:TYR:CD1	2.51	0.94
1:A:437:ASN:CA	1:A:508:TYR:CE1	2.51	0.93
1:B:604:THR:HB	5:B:1405:NAG:H82	1.50	0.93
1:C:900:MET:O	1:C:904:TYR:CD2	2.21	0.93
1:A:1107:ARG:NE	1:B:904:TYR:CD1	2.37	0.93
1:C:329:PHE:HE1	1:C:544:ASN:HA	1.29	0.93
1:C:58:PHE:HD2	1:C:290:ASP:CB	1.82	0.92
1:B:601:GLY:O	1:B:604:THR:HG23	1.68	0.92
1:C:58:PHE:CD2	1:C:290:ASP:CB	2.52	0.92
1:C:980:ILE:CD1	1:C:996:LEU:HD11	1.99	0.92
1:A:124:THR:HB	5:A:1402:NAG:H82	0.92	0.91
1:A:815:ARG:NH1	1:A:823:PHE:CE2	2.39	0.91
1:C:900:MET:HG3	1:C:904:TYR:CE2	2.06	0.91
1:A:815:ARG:NH1	1:A:823:PHE:CD2	2.39	0.91
1:A:1107:ARG:NE	1:B:904:TYR:CE1	2.39	0.91
1:A:1107:ARG:CZ	1:B:904:TYR:HD1	1.84	0.90
1:A:124:THR:CB	5:A:1402:NAG:C8	2.46	0.90
2:N:1:NAG:H62	2:N:2:NAG:C7	2.02	0.89
1:A:1107:ARG:CD	1:B:904:TYR:CD1	2.54	0.89
1:C:17:ASN:OD1	1:C:137:ASN:HB3	1.71	0.89
1:A:905:ARG:NH1	1:A:1036:GLN:HB2	1.87	0.89
1:B:601:GLY:CA	1:B:604:THR:HG22	2.03	0.89
1:B:1151:GLU:HG2	1:B:1155:TYR:HE1	1.36	0.88
1:C:618:THR:O	1:C:621:PRO:HD2	1.74	0.88
1:C:58:PHE:CE2	1:C:290:ASP:HB2	2.09	0.88
1:C:21:ARG:HG2	1:C:21:ARG:HH11	1.37	0.87
1:A:437:ASN:N	1:A:508:TYR:HD1	1.73	0.87
1:A:1107:ARG:CD	1:B:904:TYR:CE1	2.58	0.86
1:B:601:GLY:CA	1:B:604:THR:CG2	2.54	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:900:MET:O	1:A:903:ALA:HB3	1.75	0.86
1:A:1156:PHE:CZ	1:B:1155:TYR:CB	2.58	0.86
1:A:567:ARG:CD	1:B:42:VAL:HG11	2.06	0.86
1:C:329:PHE:HE1	1:C:544:ASN:CA	1.83	0.85
1:B:567:ARG:HH12	1:C:42:VAL:HG21	1.38	0.84
1:A:904:TYR:CE1	1:C:1107:ARG:NE	2.46	0.84
1:A:904:TYR:CD1	1:C:1107:ARG:CD	2.59	0.84
1:B:454:ARG:HD3	1:B:457:ARG:HG2	1.58	0.84
1:C:616:ASN:HA	1:C:644:GLN:OE1	1.77	0.84
1:C:900:MET:O	1:C:904:TYR:HD2	1.60	0.83
1:A:437:ASN:HA	1:A:508:TYR:CE1	2.10	0.83
1:A:1079:PRO:HB3	1:B:917:TYR:HE1	1.40	0.83
1:A:966:LEU:CD2	1:A:1000:ARG:HD3	2.07	0.83
1:B:966:LEU:HA	1:B:1000:ARG:HH21	1.44	0.83
1:B:1156:PHE:CE1	1:C:1155:TYR:C	2.57	0.83
1:A:34:ARG:HH21	1:A:217:PRO:HG2	1.43	0.83
1:B:1156:PHE:HE1	1:C:1155:TYR:C	1.87	0.82
1:A:437:ASN:CB	1:A:508:TYR:HE1	1.80	0.81
1:A:1079:PRO:HB2	1:B:917:TYR:CD1	2.15	0.81
1:A:904:TYR:CE1	1:C:1107:ARG:CD	2.63	0.81
1:A:904:TYR:HD1	1:C:1107:ARG:HD3	1.43	0.81
1:A:437:ASN:N	1:A:508:TYR:CD1	2.48	0.81
1:A:580:GLN:O	2:E:1:NAG:H61	1.80	0.80
1:C:273:ARG:HH12	1:C:292:ALA:CB	1.91	0.80
1:C:171:VAL:CG2	2:Z:2:NAG:H82	2.12	0.80
1:A:904:TYR:CD1	1:C:1107:ARG:NE	2.50	0.79
1:B:57:PRO:CB	1:B:273:ARG:HH22	1.94	0.78
1:B:24:LEU:HD12	1:B:25:PRO:HD2	1.65	0.78
1:A:796:ASP:OD2	1:C:709:ASN:CB	2.31	0.77
1:B:143:VAL:HG13	1:B:153:MET:O	1.84	0.77
1:A:329:PHE:CE2	1:A:528:LYS:HD3	2.19	0.77
1:B:454:ARG:HD3	1:B:457:ARG:CG	2.13	0.77
1:A:815:ARG:NH2	1:A:867:ASP:HB3	1.99	0.77
1:A:1073:LYS:HE2	1:A:1075:PHE:CZ	2.19	0.77
1:B:308:VAL:O	1:B:602:THR:HB	1.84	0.77
1:A:34:ARG:NH2	1:A:217:PRO:HG2	2.00	0.76
1:A:132:GLU:HG3	1:A:165:ASN:HB2	1.66	0.76
1:A:885:GLY:CA	1:A:901:GLN:OE1	2.33	0.76
1:A:1150:GLU:O	1:A:1154:LYS:HG3	1.85	0.76
1:A:388:ASN:HA	1:A:527:PRO:HD2	1.66	0.76
1:A:567:ARG:HG3	1:B:42:VAL:CG1	2.15	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:ARG:NH1	1:A:121:ASN:ND2	2.34	0.75
1:A:1072:GLU:OE2	1:B:894:LEU:HD23	1.86	0.75
1:A:1156:PHE:CE2	1:B:1155:TYR:HB3	2.21	0.75
1:A:521:PRO:HG2	1:B:230:PRO:O	1.87	0.74
1:C:329:PHE:HE1	1:C:544:ASN:C	1.96	0.74
1:A:332:ILE:HD12	1:A:332:ILE:N	2.02	0.74
1:C:996:LEU:O	1:C:1000:ARG:HD3	1.88	0.73
1:B:900:MET:O	1:B:904:TYR:CD2	2.41	0.73
1:B:57:PRO:HB3	1:B:273:ARG:NH2	1.99	0.73
2:c:1:NAG:H62	2:c:2:NAG:C7	2.18	0.73
1:B:517:LEU:HB2	1:C:983:ARG:NH2	2.03	0.73
1:B:604:THR:CB	5:B:1405:NAG:H82	2.17	0.73
1:A:885:GLY:HA2	1:A:901:GLN:OE1	1.89	0.73
1:C:329:PHE:CD1	1:C:544:ASN:HA	2.24	0.73
1:A:454:ARG:NH2	1:A:467:ASP:O	2.22	0.73
1:A:796:ASP:OD2	1:C:709:ASN:N	2.21	0.72
1:B:24:LEU:CD1	1:B:25:PRO:HD2	2.19	0.72
1:B:143:VAL:HG13	1:B:153:MET:C	2.14	0.72
1:A:1107:ARG:NH1	1:B:904:TYR:HD1	1.87	0.72
1:B:310:LYS:HA	1:B:600:PRO:O	1.89	0.72
1:A:897:PRO:HD2	1:A:900:MET:CE	2.19	0.72
1:A:115:GLN:HG3	1:A:132:GLU:OE1	1.90	0.72
1:A:1073:LYS:HE2	1:A:1075:PHE:HZ	1.55	0.71
1:A:815:ARG:HH21	1:A:867:ASP:CB	1.92	0.71
1:B:756:TYR:CE1	1:B:997:ILE:HG21	2.26	0.71
1:B:143:VAL:HG22	1:B:154:LYS:HG2	1.71	0.71
1:B:567:ARG:HH12	1:C:42:VAL:CG2	2.02	0.71
1:A:886:TRP:HZ3	1:A:905:ARG:HD3	1.54	0.71
1:C:273:ARG:NH1	1:C:273:ARG:HG2	2.04	0.70
1:A:1107:ARG:HD3	1:B:904:TYR:CE1	2.24	0.70
1:C:273:ARG:HG2	1:C:273:ARG:HH11	1.57	0.70
1:C:329:PHE:CE1	1:C:544:ASN:N	2.59	0.70
1:A:904:TYR:HD1	1:C:1107:ARG:CD	2.03	0.69
1:A:1152:LEU:HD21	1:B:1152:LEU:CD1	2.22	0.69
1:B:567:ARG:NH1	1:C:42:VAL:CG2	2.55	0.69
1:B:984:LEU:HD13	1:B:988:GLU:HB3	1.75	0.69
1:B:17:ASN:HB3	1:B:19:THR:HG23	1.74	0.69
1:B:1156:PHE:CE1	1:C:1155:TYR:HB3	2.27	0.68
1:A:332:ILE:HG23	1:A:362:VAL:HB	1.76	0.67
1:A:44:ARG:O	1:A:279:TYR:CD2	2.48	0.67
1:A:332:ILE:HD12	1:A:332:ILE:H	1.58	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1152:LEU:CD2	1:B:1152:LEU:HD13	2.25	0.67
1:B:144:TYR:HE2	1:B:155:SER:CB	2.08	0.67
1:B:23:GLN:OE1	1:B:23:GLN:N	2.18	0.67
1:A:886:TRP:CE3	1:A:905:ARG:CZ	2.78	0.67
1:C:544:ASN:ND2	1:C:564:GLN:HE22	1.93	0.66
1:A:1152:LEU:CD1	1:C:1152:LEU:HD11	2.24	0.66
1:A:115:GLN:CG	1:A:132:GLU:OE1	2.43	0.66
1:A:1073:LYS:HB3	1:A:1075:PHE:CE2	2.31	0.66
1:C:1000:ARG:HD2	1:C:1000:ARG:N	2.11	0.65
1:B:21:ARG:HE	1:B:78:ARG:HH11	1.42	0.65
1:B:310:LYS:CB	1:B:600:PRO:O	2.45	0.65
1:B:900:MET:O	1:B:904:TYR:HD2	1.80	0.65
1:B:396:TYR:CE2	1:C:200:TYR:HE1	2.14	0.65
1:C:357:ARG:HB2	1:C:396:TYR:CE2	2.32	0.65
1:B:310:LYS:CA	1:B:600:PRO:O	2.45	0.65
1:A:1107:ARG:NE	1:B:904:TYR:HE1	1.93	0.65
1:A:200:TYR:CD1	1:A:230:PRO:HA	2.32	0.64
1:B:900:MET:CG	1:B:904:TYR:HE2	2.02	0.64
1:C:544:ASN:ND2	1:C:564:GLN:NE2	2.45	0.64
1:C:328:ARG:NH1	1:C:580:GLN:NE2	2.45	0.64
1:B:1151:GLU:HG2	1:B:1155:TYR:CE1	2.27	0.64
1:A:1107:ARG:NE	1:B:904:TYR:HD1	1.87	0.64
1:B:457:ARG:NH1	1:B:467:ASP:OD2	2.31	0.64
1:A:1107:ARG:CD	1:B:904:TYR:HD1	2.04	0.64
2:D:1:NAG:H62	2:D:2:NAG:C7	2.28	0.64
1:C:58:PHE:HE2	1:C:275:PHE:CD2	2.16	0.63
1:A:716:THR:HG21	1:A:1073:LYS:HD3	1.80	0.63
1:A:900:MET:HE1	1:C:712:ILE:HD12	1.80	0.63
1:B:896:ILE:HD11	1:B:904:TYR:CE2	2.33	0.63
1:C:21:ARG:HH11	1:C:21:ARG:CG	2.11	0.63
1:B:520:ALA:HB1	1:B:521:PRO:HD2	1.81	0.63
1:B:1039:ARG:HD3	1:B:1042:PHE:CD2	2.34	0.63
2:Y:1:NAG:H62	2:Y:2:NAG:C7	2.28	0.63
1:A:1073:LYS:HB3	1:A:1075:PHE:HE2	1.64	0.62
1:A:200:TYR:CE1	1:A:230:PRO:HB3	2.34	0.62
1:A:896:ILE:HD11	1:A:904:TYR:CE2	2.34	0.62
1:B:52:GLN:HA	1:B:273:ARG:O	1.97	0.62
1:B:601:GLY:O	1:B:604:THR:CG2	2.43	0.62
1:A:904:TYR:HE1	1:C:1107:ARG:CD	2.12	0.62
1:B:1107:ARG:CZ	1:C:904:TYR:CD1	2.82	0.62
1:A:904:TYR:HE1	1:C:1107:ARG:HG2	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1073:LYS:CE	1:A:1075:PHE:CZ	2.83	0.62
1:A:886:TRP:HE3	1:A:905:ARG:CZ	2.13	0.61
1:B:1107:ARG:HD3	1:C:904:TYR:CD1	2.35	0.61
1:A:707:TYR:CE2	1:B:796:ASP:OD1	2.53	0.61
1:B:1150:GLU:O	1:B:1154:LYS:HG3	2.00	0.61
1:C:710:ASN:HD22	5:C:1406:NAG:H82	1.65	0.61
1:B:144:TYR:CE2	1:B:155:SER:CB	2.83	0.61
1:B:421:TYR:HB3	1:B:457:ARG:HG3	1.82	0.61
1:A:904:TYR:CE1	1:C:1107:ARG:HD3	2.30	0.61
1:A:1152:LEU:HD21	1:B:1152:LEU:HD12	1.83	0.61
1:B:149:ASN:O	1:B:150:LYS:HG2	2.01	0.60
1:A:900:MET:O	1:A:903:ALA:CB	2.49	0.60
1:B:604:THR:HA	5:B:1405:NAG:H82	1.84	0.60
1:A:896:ILE:HG13	1:A:900:MET:CE	2.31	0.60
1:C:544:ASN:HD21	1:C:564:GLN:NE2	2.00	0.60
1:A:332:ILE:CG2	1:A:362:VAL:HB	2.32	0.60
1:B:1151:GLU:O	1:B:1155:TYR:CD1	2.55	0.60
1:A:957:GLN:HG2	1:B:765:ARG:HH12	1.67	0.59
1:B:17:ASN:HB3	1:B:19:THR:CG2	2.31	0.59
1:B:980:ILE:HG21	1:B:993:ILE:CD1	2.32	0.59
1:A:332:ILE:O	1:A:332:ILE:HG22	2.02	0.59
1:B:24:LEU:HB3	1:B:25:PRO:HD2	1.84	0.59
1:A:332:ILE:HG21	1:A:362:VAL:CG1	2.32	0.59
1:A:707:TYR:HE2	1:B:796:ASP:OD1	1.84	0.59
1:C:900:MET:O	1:C:904:TYR:CE2	2.55	0.59
1:B:604:THR:CA	5:B:1405:NAG:H82	2.33	0.59
1:A:436:TRP:C	1:A:508:TYR:HD1	2.12	0.58
1:B:966:LEU:CA	1:B:1000:ARG:HH21	2.16	0.58
1:B:1156:PHE:HE1	1:C:1155:TYR:O	1.86	0.58
1:A:1152:LEU:HD13	1:C:1152:LEU:HD11	1.85	0.58
2:N:1:NAG:H62	2:N:2:NAG:O7	2.03	0.58
1:A:1017:GLU:HG2	1:B:1019:ARG:HH22	1.68	0.58
2:b:1:NAG:O4	2:b:2:NAG:O7	2.22	0.58
1:B:604:THR:HA	5:B:1405:NAG:C8	2.34	0.57
1:C:328:ARG:HH12	1:C:580:GLN:NE2	2.01	0.57
1:A:102:ARG:HH11	1:A:121:ASN:ND2	2.03	0.57
2:N:1:NAG:H62	2:N:2:NAG:C8	2.34	0.57
1:B:212:LEU:HD23	1:B:217:PRO:HB3	1.87	0.57
1:A:567:ARG:HD2	1:B:44:ARG:NH2	2.19	0.57
1:B:21:ARG:CG	1:B:78:ARG:NH1	2.68	0.57
1:A:329:PHE:CD2	1:A:528:LYS:HD3	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:748:GLU:H	1:A:748:GLU:CD	2.12	0.57
1:A:885:GLY:HA3	1:A:901:GLN:OE1	2.04	0.57
1:A:357:ARG:HG3	1:A:396:TYR:CE1	2.40	0.56
1:C:17:ASN:C	1:C:19:THR:H	2.13	0.56
1:C:329:PHE:CE1	1:C:544:ASN:C	2.81	0.56
1:C:619:GLU:O	1:C:622:VAL:HB	2.06	0.56
1:B:977:LEU:HD12	1:B:996:LEU:HD12	1.88	0.56
1:A:200:TYR:HE1	1:A:230:PRO:HB3	1.70	0.56
1:A:1107:ARG:CD	1:B:904:TYR:HE1	2.13	0.56
1:C:329:PHE:CD1	1:C:391:CYS:SG	2.99	0.56
1:A:332:ILE:HG23	1:A:362:VAL:CB	2.35	0.56
1:A:957:GLN:HG2	1:B:765:ARG:NH1	2.20	0.56
1:A:1152:LEU:HD22	1:B:1152:LEU:HD13	1.87	0.56
1:C:273:ARG:HH11	1:C:273:ARG:CG	2.17	0.56
1:B:601:GLY:C	1:B:604:THR:CG2	2.79	0.56
1:C:131:CYS:HA	1:C:165:ASN:O	2.06	0.56
1:A:331:ASN:OD1	1:A:333:THR:HG22	2.06	0.56
1:B:396:TYR:HE2	1:C:200:TYR:HE1	1.54	0.56
1:A:592:PHE:HZ	1:B:854:LYS:O	1.89	0.56
1:B:980:ILE:CG2	1:B:993:ILE:CD1	2.83	0.56
1:A:1073:LYS:CE	1:A:1075:PHE:HZ	2.18	0.55
1:B:310:LYS:HB2	1:B:600:PRO:O	2.06	0.55
1:B:1151:GLU:O	1:B:1155:TYR:HD1	1.89	0.55
1:A:340:GLU:O	1:A:344:ALA:HB2	2.07	0.55
1:A:904:TYR:HE1	1:C:1107:ARG:CG	2.20	0.55
1:B:1107:ARG:NE	1:C:904:TYR:CE1	2.74	0.55
1:A:454:ARG:HD3	1:A:457:ARG:HG2	1.88	0.55
1:A:112:SER:HB2	1:A:132:GLU:O	2.06	0.55
1:A:1072:GLU:OE2	1:B:894:LEU:CD2	2.53	0.55
1:A:716:THR:HG23	1:A:1073:LYS:HB2	1.89	0.55
1:A:896:ILE:HG13	1:A:900:MET:HE3	1.88	0.55
1:B:1101:ASP:OD2	4:V:1:NAG:H5	2.06	0.54
1:C:618:THR:C	1:C:621:PRO:HD2	2.32	0.54
1:A:712:ILE:O	1:A:1075:PHE:N	2.40	0.54
1:C:544:ASN:HD21	1:C:564:GLN:HE22	1.53	0.54
1:C:996:LEU:HB3	1:C:1000:ARG:NH2	2.23	0.54
1:A:900:MET:HG2	1:A:904:TYR:HE2	1.73	0.54
1:B:980:ILE:HG21	1:B:993:ILE:HD13	1.88	0.54
1:B:980:ILE:CG2	1:B:993:ILE:HD11	2.37	0.54
1:A:108:THR:OG1	1:A:114:THR:OG1	2.22	0.54
1:A:567:ARG:CG	1:B:42:VAL:CG1	2.86	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:966:LEU:HA	1:B:1000:ARG:NH2	2.19	0.54
1:B:900:MET:CG	1:B:904:TYR:CE2	2.83	0.53
1:C:108:THR:O	1:C:237:ARG:NE	2.41	0.53
1:B:21:ARG:HG2	1:B:78:ARG:NH1	2.23	0.53
1:C:454:ARG:CD	1:C:457:ARG:HG2	2.39	0.53
1:A:592:PHE:C	1:A:592:PHE:CD2	2.85	0.53
1:B:28:TYR:HD2	1:B:61:ASN:HB3	1.74	0.52
1:A:886:TRP:CE3	1:A:905:ARG:NH1	2.78	0.52
1:B:454:ARG:CD	1:B:457:ARG:HG2	2.35	0.52
1:B:905:ARG:NH1	1:B:1049:LEU:O	2.39	0.52
1:C:454:ARG:HD3	1:C:457:ARG:HG2	1.90	0.52
1:A:521:PRO:CG	1:B:230:PRO:O	2.55	0.52
1:B:144:TYR:HE2	1:B:155:SER:HB2	1.73	0.52
1:B:308:VAL:O	1:B:602:THR:CB	2.56	0.52
1:B:382:VAL:HG12	1:C:983:ARG:HG2	1.91	0.52
1:C:521:PRO:HA	1:C:564:GLN:OE1	2.10	0.52
1:A:904:TYR:CE1	1:C:1107:ARG:HG2	2.45	0.52
1:A:1155:TYR:HB3	1:C:1156:PHE:CZ	2.45	0.52
1:C:17:ASN:CG	1:C:137:ASN:HD22	2.18	0.52
1:B:1082:CYS:HB2	1:B:1134:ASN:OD1	2.10	0.51
1:C:329:PHE:HB3	1:C:330:PRO:HD2	1.91	0.51
1:C:905:ARG:HD2	1:C:1049:LEU:O	2.10	0.51
1:A:332:ILE:N	1:A:332:ILE:CD1	2.73	0.51
1:B:24:LEU:CD1	1:B:25:PRO:CD	2.88	0.51
1:C:21:ARG:CG	1:C:21:ARG:NH1	2.72	0.51
1:A:421:TYR:O	1:A:454:ARG:HD2	2.10	0.51
1:A:1152:LEU:HD13	1:C:1152:LEU:CD1	2.40	0.51
1:A:1094:VAL:CG2	1:B:904:TYR:OH	2.59	0.51
1:C:17:ASN:C	1:C:19:THR:N	2.69	0.51
1:A:567:ARG:CG	1:B:42:VAL:HG11	2.41	0.51
1:B:913:GLN:O	1:B:917:TYR:HD2	1.94	0.51
1:C:896:ILE:HD11	1:C:904:TYR:CE2	2.46	0.51
1:A:1073:LYS:O	1:A:1075:PHE:CD2	2.64	0.50
1:A:904:TYR:CD1	1:C:1107:ARG:CZ	2.95	0.50
1:A:1107:ARG:NH1	1:B:904:TYR:CD1	2.73	0.50
1:A:904:TYR:HE1	1:C:1107:ARG:NE	2.02	0.50
1:A:1094:VAL:HB	1:B:904:TYR:OH	2.11	0.50
1:A:1107:ARG:HG2	1:B:904:TYR:HE1	1.77	0.50
1:A:1152:LEU:HD12	1:C:1156:PHE:HE2	1.77	0.50
2:c:1:NAG:O4	2:c:2:NAG:O7	2.30	0.50
1:A:904:TYR:OH	1:C:1094:VAL:HB	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:ILE:HG21	1:A:362:VAL:HG11	1.91	0.50
1:B:1156:PHE:CE1	1:C:1156:PHE:N	2.80	0.50
1:C:127:VAL:HB	2:Z:1:NAG:O6	2.12	0.50
1:A:22:THR:O	1:A:78:ARG:NH1	2.46	0.49
1:A:708:SER:HB3	1:A:711:SER:HB3	1.93	0.49
1:A:966:LEU:HD23	1:A:1000:ARG:CD	2.33	0.49
1:B:28:TYR:CD2	1:B:61:ASN:HB3	2.46	0.49
1:B:438:SER:CB	1:B:509:ARG:HG3	2.42	0.49
1:B:213:VAL:O	1:B:213:VAL:HG12	2.12	0.49
1:B:23:GLN:H	1:B:23:GLN:CD	2.15	0.49
1:B:601:GLY:C	1:B:604:THR:HG23	2.34	0.49
1:B:644:GLN:HG2	2:R:1:NAG:H82	1.94	0.49
1:B:21:ARG:HE	1:B:78:ARG:NH1	2.10	0.49
1:B:1093:GLY:HA2	1:B:1107:ARG:HG3	1.95	0.49
1:C:17:ASN:O	1:C:19:THR:N	2.39	0.49
1:A:125:ASN:HB2	1:A:172:SER:OG	2.13	0.49
1:B:1107:ARG:NE	1:C:904:TYR:CD1	2.81	0.49
1:B:1107:ARG:CD	1:C:904:TYR:CE1	2.96	0.49
1:B:421:TYR:O	1:B:454:ARG:HD2	2.13	0.48
1:A:899:ALA:O	1:A:916:LEU:HD21	2.13	0.48
1:A:334:ASN:C	1:A:334:ASN:HD22	2.21	0.48
1:C:81:ASN:N	1:C:82:PRO:HD3	2.29	0.48
1:A:78:ARG:HH21	1:A:80:ASP:CG	2.21	0.48
1:A:404:GLY:CA	1:A:508:TYR:CD2	2.96	0.48
1:B:1152:LEU:HD11	1:B:1156:PHE:HE2	1.79	0.48
1:A:1073:LYS:O	1:A:1075:PHE:HD2	1.97	0.48
1:C:122:ASN:OD1	1:C:125:ASN:O	2.32	0.48
1:A:886:TRP:HE3	1:A:905:ARG:NH1	2.11	0.48
1:B:328:ARG:HG3	1:B:543:PHE:CE2	2.49	0.48
1:B:984:LEU:HB3	1:B:988:GLU:HB2	1.96	0.48
1:A:708:SER:HB3	1:A:711:SER:CB	2.44	0.47
1:B:147:LYS:O	1:B:149:ASN:N	2.47	0.47
1:B:421:TYR:CD2	1:B:457:ARG:HB2	2.49	0.47
1:A:900:MET:CE	1:C:712:ILE:HD12	2.44	0.47
1:A:903:ALA:HB1	1:A:913:GLN:HB2	1.95	0.47
1:A:1159:HIS:CE1	1:C:1160:THR:HG22	2.49	0.47
1:B:980:ILE:HB	1:B:993:ILE:HD11	1.96	0.47
1:C:326:ILE:HD11	1:C:534:VAL:HG22	1.96	0.47
1:C:52:GLN:HA	1:C:273:ARG:O	2.14	0.47
1:C:634:ARG:HG2	1:C:634:ARG:O	2.13	0.47
1:A:357:ARG:HH12	1:A:359:SER:HB3	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:564:GLN:HA	1:A:564:GLN:HE21	1.80	0.47
1:B:1107:ARG:HD3	1:C:904:TYR:HD1	1.78	0.47
1:A:592:PHE:CZ	1:B:854:LYS:O	2.67	0.47
1:B:149:ASN:C	1:B:151:SER:H	2.23	0.47
1:C:623:ALA:HB1	1:C:629:LEU:HD21	1.97	0.47
1:B:1107:ARG:CZ	1:C:904:TYR:HD1	2.28	0.46
1:C:21:ARG:HG2	1:C:21:ARG:NH1	2.14	0.46
1:A:567:ARG:HG3	1:B:42:VAL:HG13	1.92	0.46
1:A:1072:GLU:CD	1:B:894:LEU:CD2	2.89	0.46
1:B:29:THR:O	1:B:61:ASN:HA	2.15	0.46
1:A:708:SER:O	1:A:710:ASN:N	2.48	0.46
1:A:1072:GLU:CD	1:B:894:LEU:HD23	2.39	0.46
1:A:1152:LEU:CD2	1:B:1152:LEU:CD1	2.88	0.46
1:C:319:ARG:O	1:C:319:ARG:HG3	2.16	0.46
1:A:115:GLN:HG2	1:A:132:GLU:OE1	2.16	0.46
1:C:125:ASN:OD1	2:Z:1:NAG:H5	2.14	0.46
1:A:102:ARG:HH11	1:A:121:ASN:HD22	1.61	0.46
1:B:913:GLN:O	1:B:917:TYR:CD2	2.68	0.46
1:A:124:THR:OG1	5:A:1402:NAG:H81	2.16	0.46
1:A:708:SER:C	1:A:710:ASN:H	2.24	0.46
1:A:1107:ARG:CG	1:B:904:TYR:HE1	2.29	0.46
1:B:24:LEU:CB	1:B:25:PRO:HD2	2.45	0.46
1:C:1000:ARG:HD2	1:C:1000:ARG:H	1.80	0.46
1:B:977:LEU:HD12	1:B:996:LEU:CD1	2.46	0.46
1:B:1037:SER:OG	1:B:1039:ARG:HG3	2.16	0.46
1:A:220:PHE:CD2	1:A:220:PHE:C	2.94	0.46
1:B:143:VAL:CG2	1:B:154:LYS:HG2	2.41	0.46
1:A:567:ARG:CD	1:B:44:ARG:NH2	2.79	0.45
1:C:457:ARG:HD3	1:C:467:ASP:OD2	2.16	0.45
1:B:601:GLY:C	1:B:604:THR:HG22	2.40	0.45
1:A:332:ILE:H	1:A:332:ILE:CD1	2.27	0.45
1:A:332:ILE:CG2	1:A:362:VAL:CG1	2.94	0.45
1:B:1107:ARG:CD	1:C:904:TYR:CD1	2.99	0.45
1:B:1156:PHE:CE1	1:C:1155:TYR:O	2.67	0.45
1:C:104:TRP:HB2	1:C:106:PHE:CE1	2.51	0.45
1:B:980:ILE:CD1	1:B:996:LEU:HD11	2.47	0.45
1:B:1156:PHE:HE1	1:C:1156:PHE:N	2.13	0.45
1:B:861:LEU:H	1:B:861:LEU:HD23	1.81	0.45
1:A:132:GLU:CG	1:A:165:ASN:HB2	2.41	0.45
1:A:900:MET:O	1:A:903:ALA:N	2.49	0.45
1:B:970:PHE:O	1:B:995:ARG:HD2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:1:NAG:C6	2:N:2:NAG:C7	2.85	0.45
1:A:1107:ARG:HG2	1:B:904:TYR:CE1	2.51	0.45
1:B:164:ASN:HB3	2:P:1:NAG:H82	1.97	0.45
1:B:21:ARG:NE	1:B:78:ARG:HH11	2.13	0.45
1:B:436:TRP:CE2	1:B:509:ARG:HB2	2.52	0.45
1:A:1017:GLU:CG	1:B:1019:ARG:HH22	2.29	0.44
1:C:58:PHE:CD2	1:C:275:PHE:CE2	2.96	0.44
1:C:710:ASN:ND2	5:C:1406:NAG:H82	2.30	0.44
1:A:900:MET:HG2	1:A:904:TYR:CE2	2.50	0.44
1:A:332:ILE:CG2	1:A:362:VAL:CB	2.95	0.44
1:B:193:VAL:HG13	1:B:270:LEU:HD11	2.00	0.44
1:B:1152:LEU:HD11	1:C:1152:LEU:HD13	1.99	0.44
1:A:712:ILE:HG23	1:A:1075:PHE:HB2	2.00	0.44
1:A:1156:PHE:CE2	1:B:1155:TYR:CB	2.97	0.44
1:C:109:THR:O	1:C:237:ARG:NH1	2.51	0.44
1:C:748:GLU:CD	1:C:748:GLU:H	2.26	0.44
1:A:332:ILE:CG2	1:A:362:VAL:HG11	2.47	0.44
1:A:886:TRP:CZ3	1:A:905:ARG:CZ	3.00	0.44
1:B:454:ARG:CD	1:B:457:ARG:CG	2.92	0.44
1:B:1107:ARG:NH1	1:C:904:TYR:HD1	2.16	0.44
1:A:102:ARG:NH1	1:A:121:ASN:HD22	2.11	0.43
1:A:996:LEU:HD13	1:A:1000:ARG:NH2	2.33	0.43
1:B:64:TRP:CD1	1:B:266:TYR:CE2	3.06	0.43
1:A:437:ASN:CG	1:A:508:TYR:CE1	2.94	0.43
1:B:517:LEU:CB	1:C:983:ARG:NH2	2.79	0.43
1:B:396:TYR:CZ	1:C:230:PRO:HB3	2.53	0.43
1:B:980:ILE:HD13	1:B:993:ILE:HD13	2.01	0.43
1:A:334:ASN:C	1:A:334:ASN:ND2	2.73	0.43
1:B:168:PHE:CZ	1:B:229:LEU:HD22	2.53	0.43
1:A:804:GLN:HE22	2:H:2:NAG:H82	1.82	0.43
1:B:592:PHE:CD2	1:C:740:MET:HE1	2.54	0.43
1:B:905:ARG:HD3	1:B:1049:LEU:O	2.19	0.43
1:A:16:VAL:HG12	1:A:18:LEU:HD12	2.00	0.43
1:B:24:LEU:HB3	1:B:25:PRO:CD	2.48	0.43
1:B:1107:ARG:HD3	1:C:904:TYR:CE1	2.54	0.43
1:A:716:THR:CG2	1:A:1073:LYS:HB2	2.48	0.42
1:B:330:PRO:HG3	1:B:525:CYS:SG	2.59	0.42
1:B:1000:ARG:O	1:B:1003:SER:N	2.52	0.42
1:B:1039:ARG:HD3	1:B:1042:PHE:CG	2.53	0.42
1:B:57:PRO:CA	1:B:273:ARG:HH22	2.33	0.42
1:C:421:TYR:CD1	1:C:457:ARG:HB2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:972:ALA:HB2	1:A:995:ARG:HD2	2.02	0.42
1:A:1090:PRO:HD3	1:A:1095:PHE:CE2	2.54	0.42
1:B:1156:PHE:CZ	1:C:1155:TYR:HB3	2.55	0.42
1:A:957:GLN:CD	1:B:765:ARG:NH1	2.77	0.42
1:B:396:TYR:CE2	1:C:200:TYR:CE1	3.01	0.42
1:A:1152:LEU:CD1	1:C:1152:LEU:CD1	2.97	0.42
1:B:308:VAL:HB	1:B:602:THR:OG1	2.20	0.42
1:B:357:ARG:HG3	1:B:396:TYR:CE1	2.54	0.42
1:B:396:TYR:HE1	1:C:230:PRO:HG3	1.83	0.42
1:B:517:LEU:CB	1:C:983:ARG:HH21	2.33	0.42
1:B:756:TYR:CZ	1:B:997:ILE:CG2	3.03	0.42
1:A:1089:PHE:HB2	1:A:1121:PHE:CE2	2.55	0.42
1:B:617:CYS:N	1:B:649:CYS:SG	2.93	0.42
1:B:1094:VAL:CG2	1:C:904:TYR:OH	2.68	0.42
1:A:815:ARG:CZ	1:A:823:PHE:CE2	3.01	0.42
1:B:600:PRO:HB2	1:B:601:GLY:H	1.68	0.42
1:A:1107:ARG:CG	1:B:904:TYR:CE1	3.01	0.41
1:C:109:THR:C	1:C:237:ARG:HH11	2.28	0.41
1:A:709:ASN:HB3	1:B:796:ASP:CG	2.45	0.41
1:A:513:LEU:HD23	1:A:513:LEU:HA	1.96	0.41
1:B:330:PRO:CG	1:B:525:CYS:SG	3.08	0.41
1:B:421:TYR:HB2	1:B:454:ARG:HD2	2.02	0.41
1:C:392:PHE:CD2	1:C:517:LEU:HD21	2.55	0.41
1:C:619:GLU:O	1:C:622:VAL:N	2.54	0.41
1:B:21:ARG:CG	1:B:78:ARG:HH11	2.33	0.41
1:C:1000:ARG:N	1:C:1000:ARG:CD	2.81	0.41
1:B:104:TRP:HB2	1:B:106:PHE:CE1	2.56	0.40
1:B:604:THR:HB	5:B:1405:NAG:C8	2.36	0.40
1:A:957:GLN:CG	1:B:765:ARG:NH1	2.84	0.40
1:B:21:ARG:HG2	1:B:78:ARG:HH11	1.86	0.40
1:B:1039:ARG:NE	1:C:1031:GLU:OE2	2.49	0.40
1:A:124:THR:CB	5:A:1402:NAG:H81	2.43	0.40
1:B:445:VAL:HA	1:B:499:PRO:HD2	2.03	0.40
1:B:1039:ARG:HG3	1:B:1039:ARG:H	1.74	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	954/1310 (73%)	874 (92%)	71 (7%)	9 (1%)	14	49
1	B	1047/1310 (80%)	951 (91%)	89 (8%)	7 (1%)	18	55
1	C	1083/1310 (83%)	1001 (92%)	78 (7%)	4 (0%)	30	66
All	All	3084/3930 (78%)	2826 (92%)	238 (8%)	20 (1%)	23	58

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	ASN
1	A	709	ASN
1	B	146	HIS
1	B	148	ASN
1	B	600	PRO
1	C	291	CYS
1	A	148	ASN
1	A	393	THR
1	A	737	ASP
1	A	758	SER
1	A	834	ILE
1	B	620	VAL
1	A	381	GLY
1	C	353	TRP
1	A	603	ASN
1	B	619	GLU
1	C	20	THR
1	C	1126	CYS
1	B	812	PRO
1	B	257	GLY

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	448/1135 (40%)	432 (96%)	16 (4%)	31	52
1	B	21/1135 (2%)	21 (100%)	0	100	100
All	All	469/2270 (21%)	453 (97%)	16 (3%)	33	54

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	132	GLU
1	A	213	VAL
1	A	220	PHE
1	A	266	TYR
1	A	282	ASN
1	A	332	ILE
1	A	333	THR
1	A	335	LEU
1	A	346	ARG
1	A	351	TYR
1	A	367	VAL
1	A	368	LEU
1	A	392	PHE
1	A	461	LEU
1	A	764	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	ASN
1	A	121	ASN
1	A	137	ASN
1	A	188	ASN
1	A	196	ASN
1	A	334	ASN

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Mol	Chain	Res	Type
1	A	354	ASN
1	A	370	ASN
1	A	450	ASN
1	A	460	ASN
1	A	644	GLN
1	A	655	HIS
1	A	762	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 ⓘ

80 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	D	1	2,1	14,14,15	0.30	0	17,19,21	0.39	0
2	NAG	D	2	2	14,14,15	1.15	1 (7%)	17,19,21	0.92	1 (5%)
2	NAG	E	1	2,1	14,14,15	1.41	2 (14%)	17,19,21	1.71	2 (11%)
2	NAG	E	2	2	14,14,15	1.20	1 (7%)	17,19,21	0.63	0
2	NAG	F	1	2,1	14,14,15	1.24	1 (7%)	17,19,21	1.22	3 (17%)
2	NAG	F	2	2	14,14,15	1.25	2 (14%)	17,19,21	0.76	0
2	NAG	G	1	2,1	14,14,15	1.16	1 (7%)	17,19,21	1.21	2 (11%)
2	NAG	G	2	2	14,14,15	1.20	1 (7%)	17,19,21	0.71	0
2	NAG	H	1	2,1	14,14,15	1.13	1 (7%)	17,19,21	0.74	0
2	NAG	H	2	2	14,14,15	1.20	1 (7%)	17,19,21	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	I	1	3,1	14,14,15	1.45	3 (21%)	17,19,21	1.29	2 (11%)
3	NAG	I	2	3	14,14,15	1.24	1 (7%)	17,19,21	0.92	1 (5%)
3	FUC	I	3	3	10,10,11	1.53	2 (20%)	14,14,16	1.00	0
4	NAG	J	1	1,4	14,14,15	1.22	1 (7%)	17,19,21	1.14	2 (11%)
4	NAG	J	2	4	14,14,15	1.38	3 (21%)	17,19,21	0.46	0
4	MAN	J	3	4	11,11,12	1.49	2 (18%)	15,15,17	1.02	1 (6%)
4	NAG	K	1	1,4	14,14,15	0.23	0	17,19,21	0.53	0
4	NAG	K	2	4	14,14,15	1.26	2 (14%)	17,19,21	0.57	0
4	MAN	K	3	4	11,11,12	1.48	2 (18%)	15,15,17	0.89	0
4	NAG	L	1	1,4	14,14,15	1.53	3 (21%)	17,19,21	1.15	1 (5%)
4	NAG	L	2	4	14,14,15	1.39	2 (14%)	17,19,21	0.80	0
4	MAN	L	3	4	11,11,12	1.42	1 (9%)	15,15,17	0.98	2 (13%)
4	NAG	M	1	1,4	14,14,15	1.28	1 (7%)	17,19,21	1.33	1 (5%)
4	NAG	M	2	4	14,14,15	1.38	2 (14%)	17,19,21	0.73	0
4	MAN	M	3	4	11,11,12	1.50	2 (18%)	15,15,17	0.88	0
2	NAG	N	1	2,1	14,14,15	0.29	0	17,19,21	0.38	0
2	NAG	N	2	2	14,14,15	1.10	1 (7%)	17,19,21	0.72	0
2	NAG	O	1	2,1	14,14,15	1.16	2 (14%)	17,19,21	0.95	1 (5%)
2	NAG	O	2	2	14,14,15	1.27	3 (21%)	17,19,21	1.26	1 (5%)
2	NAG	P	1	2,1	14,14,15	0.25	0	17,19,21	0.53	0
2	NAG	P	2	2	14,14,15	0.28	0	17,19,21	0.55	0
2	NAG	Q	1	2,1	14,14,15	1.22	1 (7%)	17,19,21	0.88	0
2	NAG	Q	2	2	14,14,15	1.50	3 (21%)	17,19,21	0.90	1 (5%)
2	NAG	R	1	2,1	14,14,15	0.29	0	17,19,21	0.45	0
2	NAG	R	2	2	14,14,15	1.19	1 (7%)	17,19,21	0.81	1 (5%)
2	NAG	S	1	2,1	14,14,15	1.06	1 (7%)	17,19,21	1.00	0
2	NAG	S	2	2	14,14,15	1.13	1 (7%)	17,19,21	0.66	0
2	NAG	T	1	2,1	14,14,15	1.16	2 (14%)	17,19,21	1.04	1 (5%)
2	NAG	T	2	2	14,14,15	1.41	2 (14%)	17,19,21	0.83	1 (5%)
3	NAG	U	1	3,1	14,14,15	0.29	0	17,19,21	0.92	1 (5%)
3	NAG	U	2	3	14,14,15	0.29	0	17,19,21	0.55	0
3	FUC	U	3	3	10,10,11	1.53	1 (10%)	14,14,16	1.90	3 (21%)
4	NAG	V	1	1,4	14,14,15	1.52	3 (21%)	17,19,21	1.18	2 (11%)
4	NAG	V	2	4	14,14,15	1.35	1 (7%)	17,19,21	1.53	3 (17%)
4	MAN	V	3	4	11,11,12	1.43	2 (18%)	15,15,17	0.61	0
4	NAG	W	1	1,4	14,14,15	0.22	0	17,19,21	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	W	2	4	14,14,15	1.27	2 (14%)	17,19,21	0.56	0
4	MAN	W	3	4	11,11,12	1.50	2 (18%)	15,15,17	0.88	0
4	NAG	X	1	1,4	14,14,15	1.29	1 (7%)	17,19,21	1.33	1 (5%)
4	NAG	X	2	4	14,14,15	1.38	2 (14%)	17,19,21	0.73	0
4	MAN	X	3	4	11,11,12	1.51	2 (18%)	15,15,17	0.89	0
2	NAG	Y	1	2,1	14,14,15	0.29	0	17,19,21	0.39	0
2	NAG	Y	2	2	14,14,15	1.16	1 (7%)	17,19,21	0.92	1 (5%)
2	NAG	Z	1	2,1	14,14,15	0.21	0	17,19,21	0.41	0
2	NAG	Z	2	2	14,14,15	1.06	1 (7%)	17,19,21	1.42	1 (5%)
2	NAG	a	1	2,1	14,14,15	1.27	1 (7%)	17,19,21	0.75	0
2	NAG	a	2	2	14,14,15	1.25	1 (7%)	17,19,21	1.23	2 (11%)
2	NAG	b	1	2,1	14,14,15	1.31	1 (7%)	17,19,21	1.46	3 (17%)
2	NAG	b	2	2	14,14,15	1.24	1 (7%)	17,19,21	1.03	1 (5%)
2	NAG	c	1	2,1	14,14,15	0.27	0	17,19,21	0.46	0
2	NAG	c	2	2	14,14,15	1.06	1 (7%)	17,19,21	1.32	1 (5%)
2	NAG	d	1	2,1	14,14,15	1.13	1 (7%)	17,19,21	0.87	0
2	NAG	d	2	2	14,14,15	1.05	1 (7%)	17,19,21	0.81	0
2	NAG	e	1	2,1	14,14,15	1.27	2 (14%)	17,19,21	0.92	0
2	NAG	e	2	2	14,14,15	1.27	3 (21%)	17,19,21	0.84	0
3	NAG	f	1	3,1	14,14,15	1.37	3 (21%)	17,19,21	1.02	1 (5%)
3	NAG	f	2	3	14,14,15	1.27	2 (14%)	17,19,21	0.80	0
3	FUC	f	3	3	10,10,11	1.71	2 (20%)	14,14,16	1.18	2 (14%)
4	NAG	g	1	1,4	14,14,15	1.14	2 (14%)	17,19,21	1.07	1 (5%)
4	NAG	g	2	4	14,14,15	1.38	3 (21%)	17,19,21	1.22	1 (5%)
4	MAN	g	3	4	11,11,12	1.46	2 (18%)	15,15,17	0.96	1 (6%)
4	NAG	h	1	1,4	14,14,15	1.44	3 (21%)	17,19,21	0.96	0
4	NAG	h	2	4	14,14,15	1.39	3 (21%)	17,19,21	0.99	2 (11%)
4	MAN	h	3	4	11,11,12	1.59	2 (18%)	15,15,17	0.93	1 (6%)
4	NAG	i	1	1,4	14,14,15	0.28	0	17,19,21	0.58	0
4	NAG	i	2	4	14,14,15	0.28	0	17,19,21	0.54	0
4	MAN	i	3	4	11,11,12	0.23	0	15,15,17	0.50	0
4	NAG	j	1	1,4	14,14,15	1.53	3 (21%)	17,19,21	1.17	2 (11%)
4	NAG	j	2	4	14,14,15	1.39	2 (14%)	17,19,21	0.62	0
4	MAN	j	3	4	11,11,12	1.56	2 (18%)	15,15,17	0.90	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

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

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	U	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	U	2	3	-	0/6/23/26	0/1/1/1
3	FUC	U	3	3	-	-	0/1/1/1
4	NAG	V	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	V	2	4	-	0/6/23/26	0/1/1/1
4	MAN	V	3	4	-	1/2/19/22	0/1/1/1
4	NAG	W	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	W	2	4	-	0/6/23/26	0/1/1/1
4	MAN	W	3	4	-	0/2/19/22	0/1/1/1
4	NAG	X	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	X	2	4	-	0/6/23/26	0/1/1/1
4	MAN	X	3	4	-	1/2/19/22	0/1/1/1
2	NAG	Y	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	Y	2	2	-	0/6/23/26	0/1/1/1
2	NAG	Z	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	Z	2	2	-	0/6/23/26	0/1/1/1
2	NAG	a	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	a	2	2	-	0/6/23/26	0/1/1/1
2	NAG	b	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	b	2	2	-	1/6/23/26	0/1/1/1
2	NAG	c	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	c	2	2	-	1/6/23/26	0/1/1/1
2	NAG	d	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	d	2	2	-	0/6/23/26	0/1/1/1
2	NAG	e	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	e	2	2	-	0/6/23/26	0/1/1/1
3	NAG	f	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	f	2	3	-	0/6/23/26	0/1/1/1
3	FUC	f	3	3	-	-	0/1/1/1
4	NAG	g	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	g	2	4	-	0/6/23/26	0/1/1/1
4	MAN	g	3	4	-	1/2/19/22	1/1/1/1
4	NAG	h	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	h	2	4	-	0/6/23/26	0/1/1/1
4	MAN	h	3	4	-	0/2/19/22	1/1/1/1
4	NAG	i	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	i	2	4	-	1/6/23/26	0/1/1/1
4	MAN	i	3	4	-	1/2/19/22	1/1/1/1
4	NAG	j	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	j	2	4	-	0/6/23/26	0/1/1/1
4	MAN	j	3	4	-	1/2/19/22	0/1/1/1

All (114) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1	NAG	O5-C5	3.63	1.50	1.43
4	h	3	MAN	O5-C5	3.40	1.50	1.43
4	j	1	NAG	C1-C2	3.33	1.56	1.52
4	V	2	NAG	O5-C5	3.31	1.49	1.43
4	g	3	MAN	O5-C5	3.31	1.49	1.43
2	Q	2	NAG	O5-C5	3.30	1.49	1.43
4	V	1	NAG	C1-C2	3.27	1.56	1.52
4	L	1	NAG	C1-C2	3.25	1.56	1.52
2	R	2	NAG	O5-C5	3.20	1.49	1.43
2	e	1	NAG	O5-C5	3.16	1.49	1.43
4	h	1	NAG	O5-C5	3.15	1.49	1.43
2	Y	2	NAG	O5-C5	3.13	1.49	1.43
4	W	3	MAN	O5-C5	3.13	1.49	1.43
4	j	3	MAN	O5-C5	3.12	1.49	1.43
2	D	2	NAG	O5-C5	3.12	1.49	1.43
4	L	2	NAG	O5-C5	3.09	1.49	1.43
2	T	2	NAG	O5-C1	3.06	1.48	1.43
4	K	3	MAN	O5-C5	3.06	1.49	1.43
4	j	2	NAG	O5-C5	3.01	1.49	1.43
3	U	3	FUC	O5-C5	3.00	1.49	1.43
4	L	3	MAN	O5-C5	2.92	1.49	1.43
4	J	3	MAN	O5-C5	2.92	1.49	1.43
4	X	1	NAG	O5-C5	2.89	1.49	1.43
2	b	1	NAG	O5-C5	2.88	1.49	1.43
4	M	1	NAG	O5-C5	2.87	1.49	1.43
2	F	1	NAG	O5-C5	2.86	1.49	1.43
2	G	2	NAG	O5-C5	2.86	1.49	1.43
3	f	3	FUC	O5-C5	2.86	1.49	1.43
4	V	3	MAN	O5-C5	2.85	1.49	1.43
3	I	1	NAG	O5-C5	2.84	1.49	1.43
2	a	2	NAG	O5-C5	2.83	1.49	1.43
3	f	3	FUC	O5-C1	2.82	1.48	1.43
4	W	2	NAG	O5-C5	2.82	1.48	1.43
2	H	2	NAG	O5-C5	2.82	1.48	1.43
2	Q	2	NAG	O5-C1	2.80	1.48	1.43
4	J	2	NAG	O5-C5	2.78	1.48	1.43
4	K	2	NAG	O5-C5	2.78	1.48	1.43
3	I	2	NAG	O5-C5	2.78	1.48	1.43
4	X	3	MAN	O5-C5	2.76	1.48	1.43
2	T	1	NAG	O5-C5	2.75	1.48	1.43
3	I	1	NAG	C1-C2	2.74	1.56	1.52
2	F	2	NAG	O5-C5	2.74	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	J	1	NAG	O5-C5	2.72	1.48	1.43
4	M	3	MAN	O5-C5	2.72	1.48	1.43
2	O	2	NAG	O5-C5	2.69	1.48	1.43
2	T	2	NAG	O5-C5	2.68	1.48	1.43
3	I	3	FUC	O5-C5	2.67	1.48	1.43
2	b	2	NAG	O5-C5	2.66	1.48	1.43
4	j	2	NAG	O4-C4	2.65	1.49	1.43
2	E	2	NAG	O5-C5	2.65	1.48	1.43
4	K	2	NAG	O4-C4	2.64	1.49	1.43
2	c	2	NAG	O5-C5	2.63	1.48	1.43
2	a	1	NAG	O5-C5	2.63	1.48	1.43
4	h	2	NAG	O5-C5	2.62	1.48	1.43
4	W	2	NAG	O4-C4	2.61	1.49	1.43
4	J	3	MAN	O5-C1	2.60	1.48	1.43
4	X	2	NAG	O4-C4	2.55	1.49	1.43
2	e	2	NAG	O5-C5	2.55	1.48	1.43
4	M	2	NAG	O4-C4	2.54	1.49	1.43
3	f	1	NAG	C1-C2	2.53	1.55	1.52
3	f	1	NAG	O5-C5	2.52	1.48	1.43
2	Z	2	NAG	O5-C5	2.52	1.48	1.43
4	X	2	NAG	O5-C5	2.51	1.48	1.43
2	G	1	NAG	O5-C5	2.50	1.48	1.43
3	f	2	NAG	O5-C5	2.50	1.48	1.43
4	M	2	NAG	O5-C5	2.49	1.48	1.43
4	g	2	NAG	O5-C1	2.43	1.47	1.43
4	L	2	NAG	O4-C4	2.43	1.49	1.43
2	Q	1	NAG	O5-C5	2.42	1.48	1.43
4	h	3	MAN	O5-C1	2.41	1.47	1.43
4	V	1	NAG	O5-C5	2.41	1.48	1.43
4	L	1	NAG	O5-C5	2.40	1.48	1.43
4	j	1	NAG	O5-C5	2.39	1.48	1.43
4	h	1	NAG	O5-C1	2.38	1.47	1.43
3	I	3	FUC	O5-C1	2.36	1.47	1.43
2	S	2	NAG	O5-C5	2.35	1.48	1.43
2	N	2	NAG	O5-C5	2.33	1.48	1.43
2	O	1	NAG	O5-C5	2.33	1.48	1.43
4	K	3	MAN	O5-C1	2.31	1.47	1.43
4	g	1	NAG	O5-C1	2.30	1.47	1.43
2	d	2	NAG	O5-C5	2.29	1.47	1.43
4	W	3	MAN	O5-C1	2.29	1.47	1.43
2	H	1	NAG	O5-C5	2.27	1.47	1.43
2	e	2	NAG	C1-C2	2.25	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	d	1	NAG	C1-C2	2.24	1.55	1.52
4	j	1	NAG	O4-C4	2.24	1.48	1.43
4	g	3	MAN	O5-C1	2.23	1.47	1.43
3	f	1	NAG	O4-C4	2.23	1.48	1.43
4	L	1	NAG	O4-C4	2.22	1.48	1.43
4	h	2	NAG	O4-C4	2.21	1.48	1.43
4	h	2	NAG	O5-C1	2.18	1.47	1.43
4	V	1	NAG	O4-C4	2.17	1.48	1.43
4	h	1	NAG	C1-C2	2.16	1.55	1.52
4	J	2	NAG	O4-C4	2.15	1.48	1.43
4	V	3	MAN	O5-C1	2.15	1.47	1.43
4	g	1	NAG	O4-C4	2.14	1.48	1.43
4	J	2	NAG	O5-C1	2.12	1.47	1.43
4	X	3	MAN	O5-C1	2.12	1.47	1.43
4	M	3	MAN	O5-C1	2.10	1.47	1.43
2	e	1	NAG	C8-C7	2.10	1.54	1.50
3	f	2	NAG	C1-C2	2.08	1.55	1.52
2	Q	2	NAG	C1-C2	2.07	1.55	1.52
2	S	1	NAG	O5-C5	2.07	1.47	1.43
2	O	2	NAG	O5-C1	2.07	1.47	1.43
4	j	3	MAN	O5-C1	2.07	1.47	1.43
4	g	2	NAG	O5-C5	2.06	1.47	1.43
2	O	2	NAG	C1-C2	2.04	1.55	1.52
4	g	2	NAG	C1-C2	2.03	1.55	1.52
2	T	1	NAG	C8-C7	2.01	1.54	1.50
2	e	2	NAG	O5-C1	2.01	1.47	1.43
2	F	2	NAG	O5-C1	2.01	1.47	1.43
3	I	1	NAG	O4-C4	2.00	1.47	1.43
2	O	1	NAG	O4-C4	2.00	1.47	1.43
2	E	1	NAG	O4-C4	2.00	1.47	1.43

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	NAG	C1-O5-C5	5.69	119.82	112.19
3	U	3	FUC	C1-C2-C3	5.02	116.95	109.64
2	Z	2	NAG	C1-O5-C5	4.97	118.84	112.19
2	O	2	NAG	C1-O5-C5	4.76	118.56	112.19
2	c	2	NAG	C1-O5-C5	4.64	118.40	112.19
4	V	2	NAG	C1-O5-C5	4.23	117.85	112.19
3	I	1	NAG	C1-O5-C5	4.16	117.76	112.19
4	X	1	NAG	C1-O5-C5	4.12	117.70	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	M	1	NAG	C1-O5-C5	4.11	117.69	112.19
4	g	2	NAG	C1-O5-C5	4.07	117.64	112.19
2	b	1	NAG	C3-C4-C5	3.62	116.80	110.23
2	a	2	NAG	C1-O5-C5	3.31	116.62	112.19
3	U	3	FUC	C1-O5-C5	3.29	120.74	112.97
2	E	1	NAG	C4-C3-C2	-3.22	106.30	111.02
4	g	1	NAG	C1-O5-C5	3.14	116.39	112.19
3	U	1	NAG	C1-O5-C5	3.12	116.36	112.19
2	b	1	NAG	C1-O5-C5	3.11	116.36	112.19
2	b	2	NAG	C1-O5-C5	2.96	116.16	112.19
3	f	1	NAG	C1-O5-C5	2.93	116.12	112.19
4	J	1	NAG	C1-O5-C5	2.85	116.00	112.19
2	G	1	NAG	C1-O5-C5	2.84	115.99	112.19
4	V	1	NAG	O5-C1-C2	-2.80	106.95	111.29
4	j	1	NAG	O5-C1-C2	-2.76	107.02	111.29
4	L	1	NAG	O5-C1-C2	-2.70	107.11	111.29
2	F	1	NAG	C4-C3-C2	-2.64	107.14	111.02
2	R	2	NAG	C1-O5-C5	2.63	115.71	112.19
3	U	3	FUC	C3-C4-C5	-2.62	105.83	109.81
4	h	2	NAG	O5-C1-C2	-2.56	107.33	111.29
3	I	1	NAG	O5-C5-C6	2.56	112.65	107.66
2	F	1	NAG	O5-C1-C2	-2.49	107.44	111.29
2	G	1	NAG	C4-C3-C2	-2.44	107.44	111.02
4	L	3	MAN	C1-O5-C5	2.42	115.42	112.19
4	J	3	MAN	C1-O5-C5	2.38	115.38	112.19
4	J	1	NAG	O4-C4-C3	-2.34	104.86	110.38
3	I	2	NAG	C1-O5-C5	2.28	115.24	112.19
2	O	1	NAG	O5-C1-C2	-2.27	107.78	111.29
2	Q	2	NAG	C1-O5-C5	2.25	115.20	112.19
2	a	2	NAG	C3-C4-C5	2.24	114.30	110.23
2	b	1	NAG	O5-C1-C2	-2.20	107.89	111.29
2	T	2	NAG	C1-O5-C5	2.19	115.12	112.19
2	Y	2	NAG	O4-C4-C3	-2.17	105.26	110.38
4	V	2	NAG	C6-C5-C4	-2.16	107.72	113.02
2	D	2	NAG	O4-C4-C3	-2.15	105.30	110.38
4	V	2	NAG	C3-C4-C5	2.14	114.11	110.23
4	g	3	MAN	C1-C2-C3	2.14	112.76	109.64
4	h	3	MAN	C1-C2-C3	2.12	112.73	109.64
4	L	3	MAN	C1-C2-C3	2.11	112.72	109.64
4	V	1	NAG	C1-O5-C5	2.10	115.00	112.19
2	F	1	NAG	C1-O5-C5	2.09	114.98	112.19
2	T	1	NAG	C1-O5-C5	-2.08	109.39	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	j	1	NAG	C1-O5-C5	2.06	114.95	112.19
3	f	3	FUC	C1-O5-C5	2.05	117.79	112.97
4	j	3	MAN	C1-O5-C5	2.02	114.89	112.19
4	h	2	NAG	C1-O5-C5	2.02	114.89	112.19
3	f	3	FUC	O2-C2-C1	2.02	113.85	109.22

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	f	1	NAG	O5-C5-C6-O6
3	U	1	NAG	O5-C5-C6-O6
3	I	1	NAG	O5-C5-C6-O6
3	f	1	NAG	C4-C5-C6-O6
3	I	1	NAG	C4-C5-C6-O6
2	b	1	NAG	C4-C5-C6-O6
3	U	1	NAG	C4-C5-C6-O6
4	K	1	NAG	O5-C5-C6-O6
4	W	1	NAG	O5-C5-C6-O6
4	K	1	NAG	C4-C5-C6-O6
4	W	1	NAG	C4-C5-C6-O6
2	N	1	NAG	O5-C5-C6-O6
2	b	1	NAG	O5-C5-C6-O6
4	j	3	MAN	O5-C5-C6-O6
4	J	3	MAN	O5-C5-C6-O6
4	M	3	MAN	O5-C5-C6-O6
4	X	3	MAN	O5-C5-C6-O6
4	L	3	MAN	O5-C5-C6-O6
2	H	1	NAG	O5-C5-C6-O6
4	g	3	MAN	O5-C5-C6-O6
4	i	3	MAN	O5-C5-C6-O6
4	i	2	NAG	O5-C5-C6-O6
4	V	3	MAN	O5-C5-C6-O6
4	i	1	NAG	C8-C7-N2-C2
4	i	1	NAG	O7-C7-N2-C2
2	N	1	NAG	C4-C5-C6-O6
2	b	2	NAG	C1-C2-N2-C7
2	E	1	NAG	C4-C5-C6-O6
2	H	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C8-C7-N2-C2
2	S	1	NAG	O5-C5-C6-O6
2	c	2	NAG	C4-C5-C6-O6

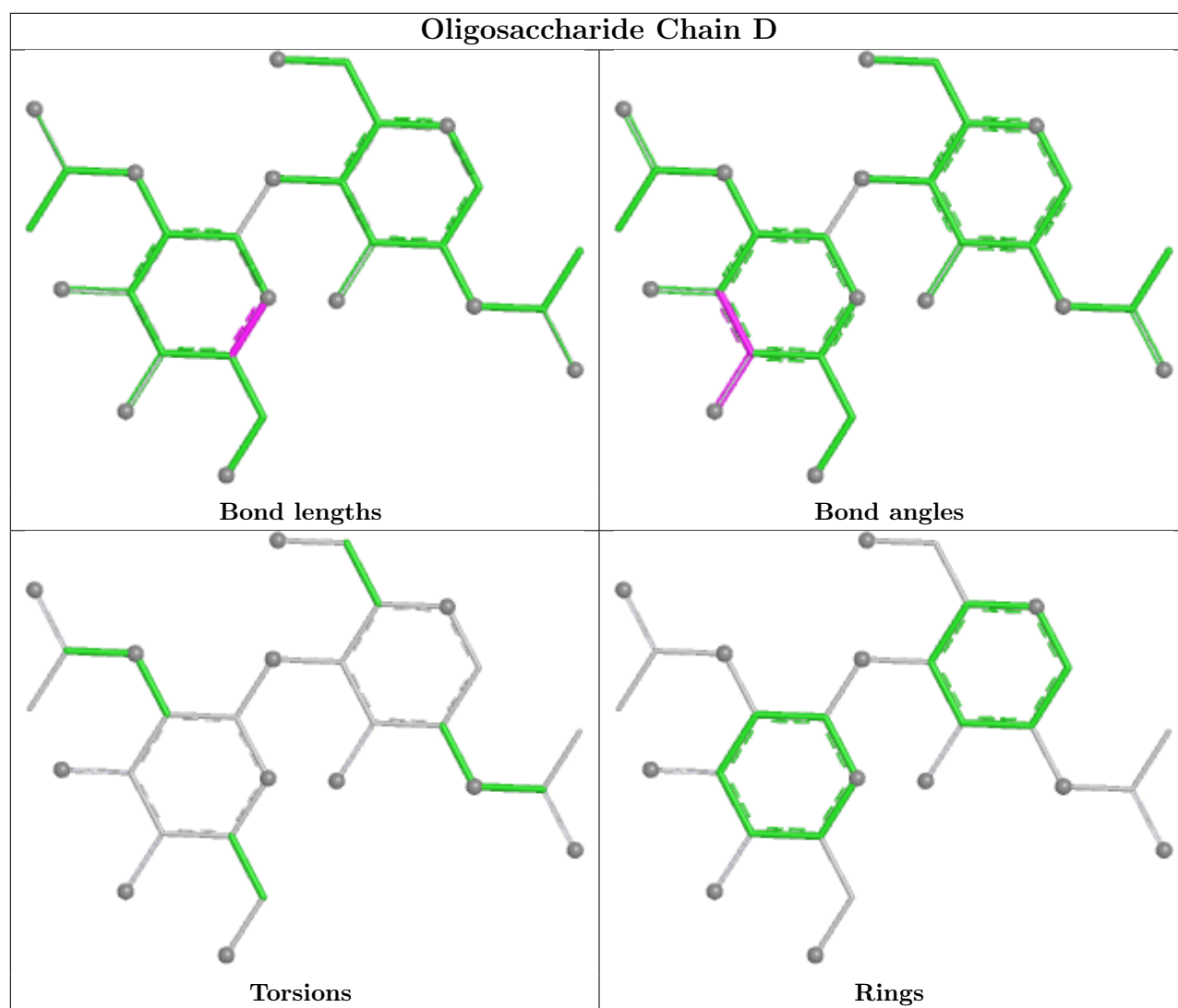
All (4) ring outliers are listed below:

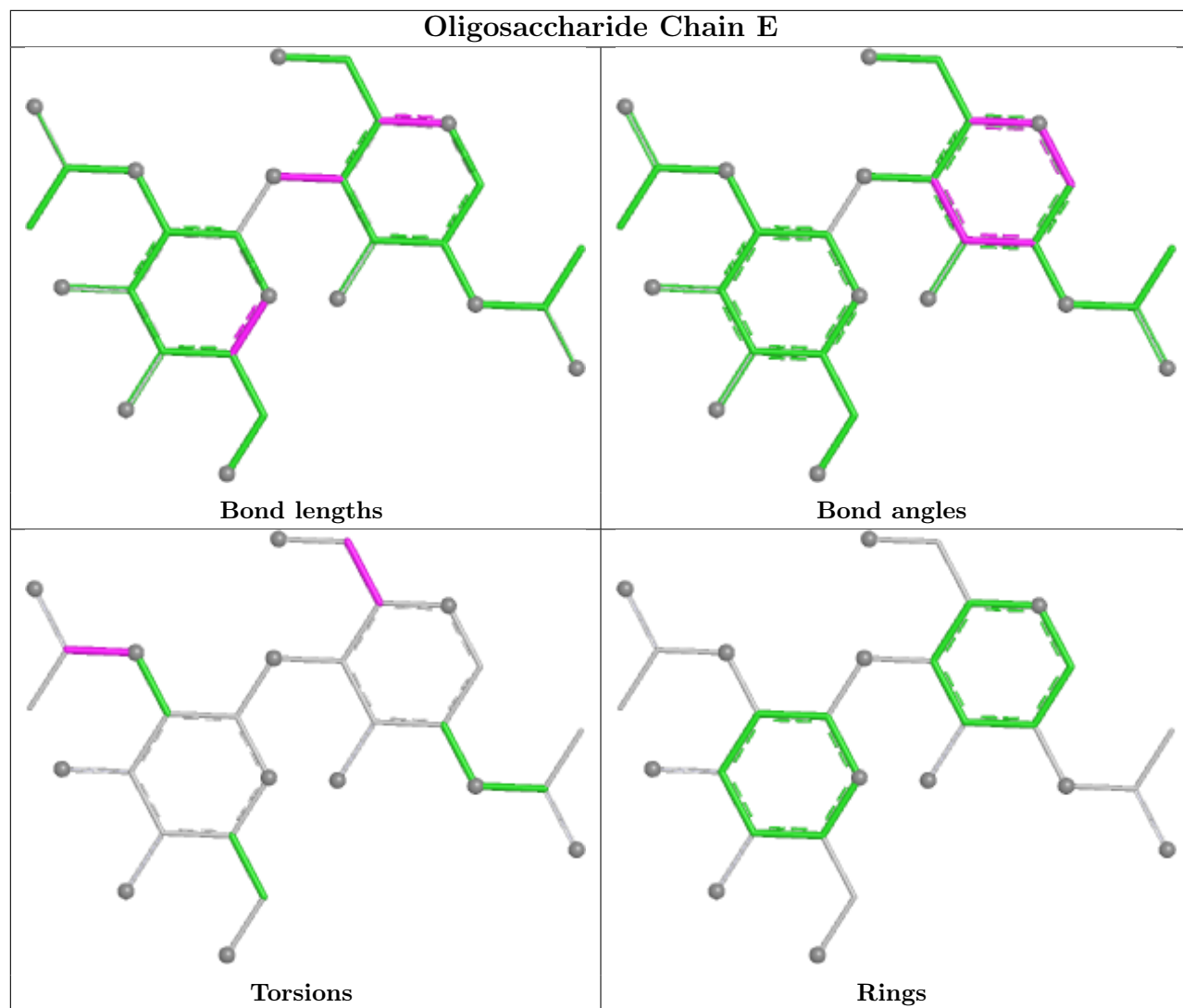
Mol	Chain	Res	Type	Atoms
4	h	3	MAN	C1-C2-C3-C4-C5-O5
4	i	3	MAN	C1-C2-C3-C4-C5-O5
4	L	3	MAN	C1-C2-C3-C4-C5-O5
4	g	3	MAN	C1-C2-C3-C4-C5-O5

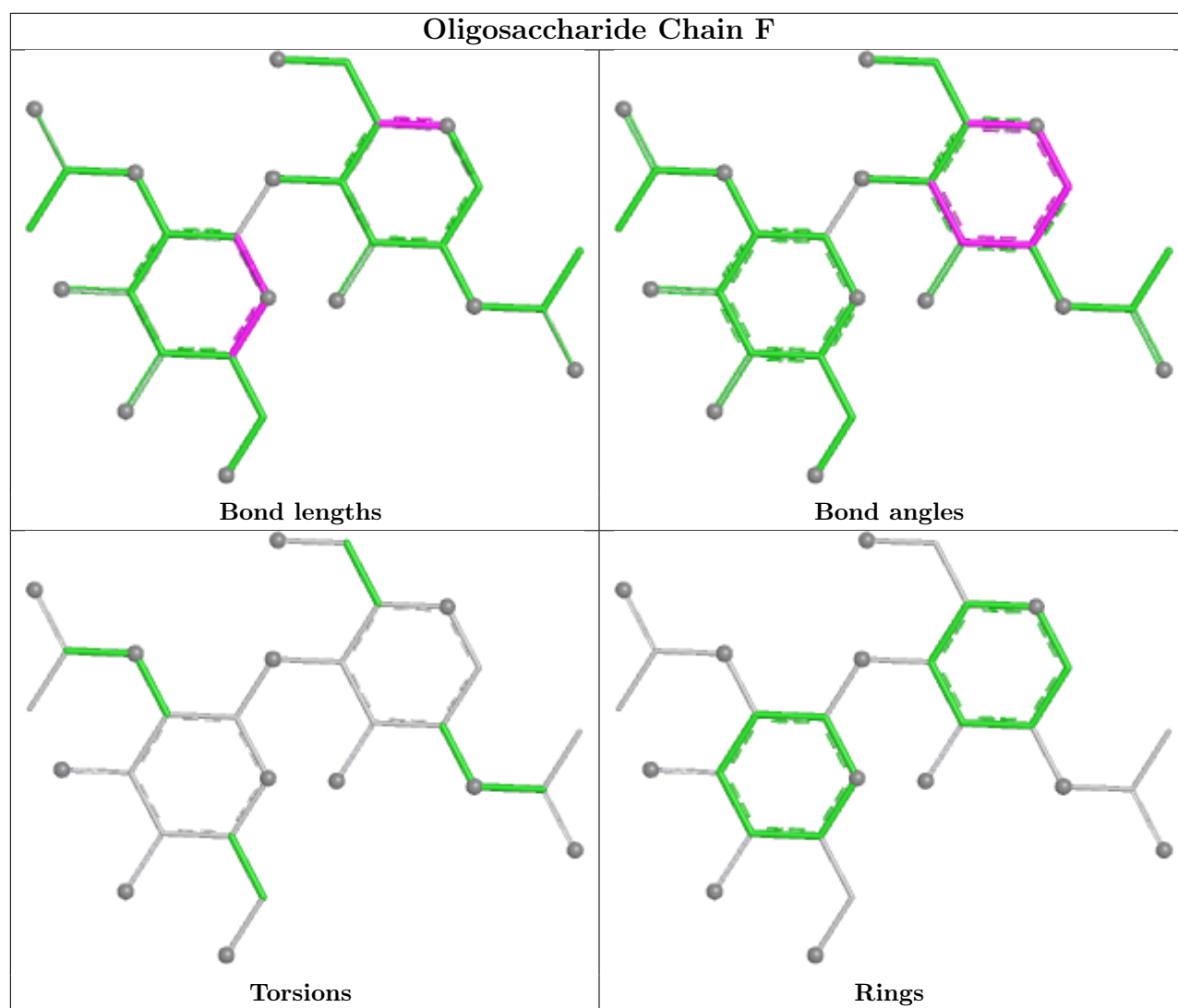
17 monomers are involved in 18 short contacts:

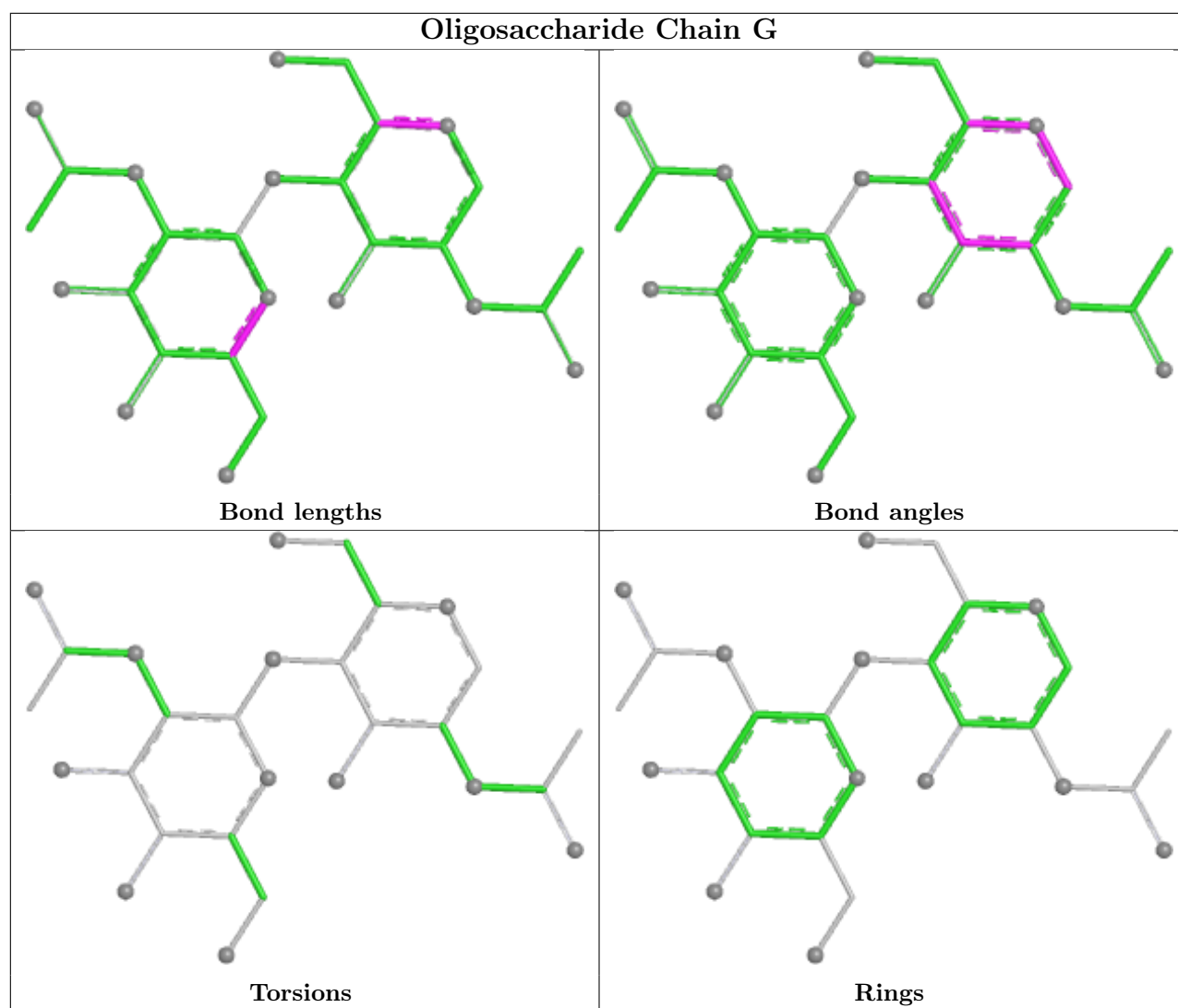
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	N	1	NAG	4	0
2	D	1	NAG	1	0
2	c	2	NAG	2	0
2	c	1	NAG	2	0
2	P	1	NAG	1	0
2	b	1	NAG	1	0
2	D	2	NAG	1	0
2	R	1	NAG	1	0
2	Z	2	NAG	2	0
2	Y	2	NAG	1	0
2	E	1	NAG	1	0
2	b	2	NAG	1	0
2	H	2	NAG	1	0
2	Y	1	NAG	1	0
4	V	1	NAG	1	0
2	N	2	NAG	4	0
2	Z	1	NAG	2	0

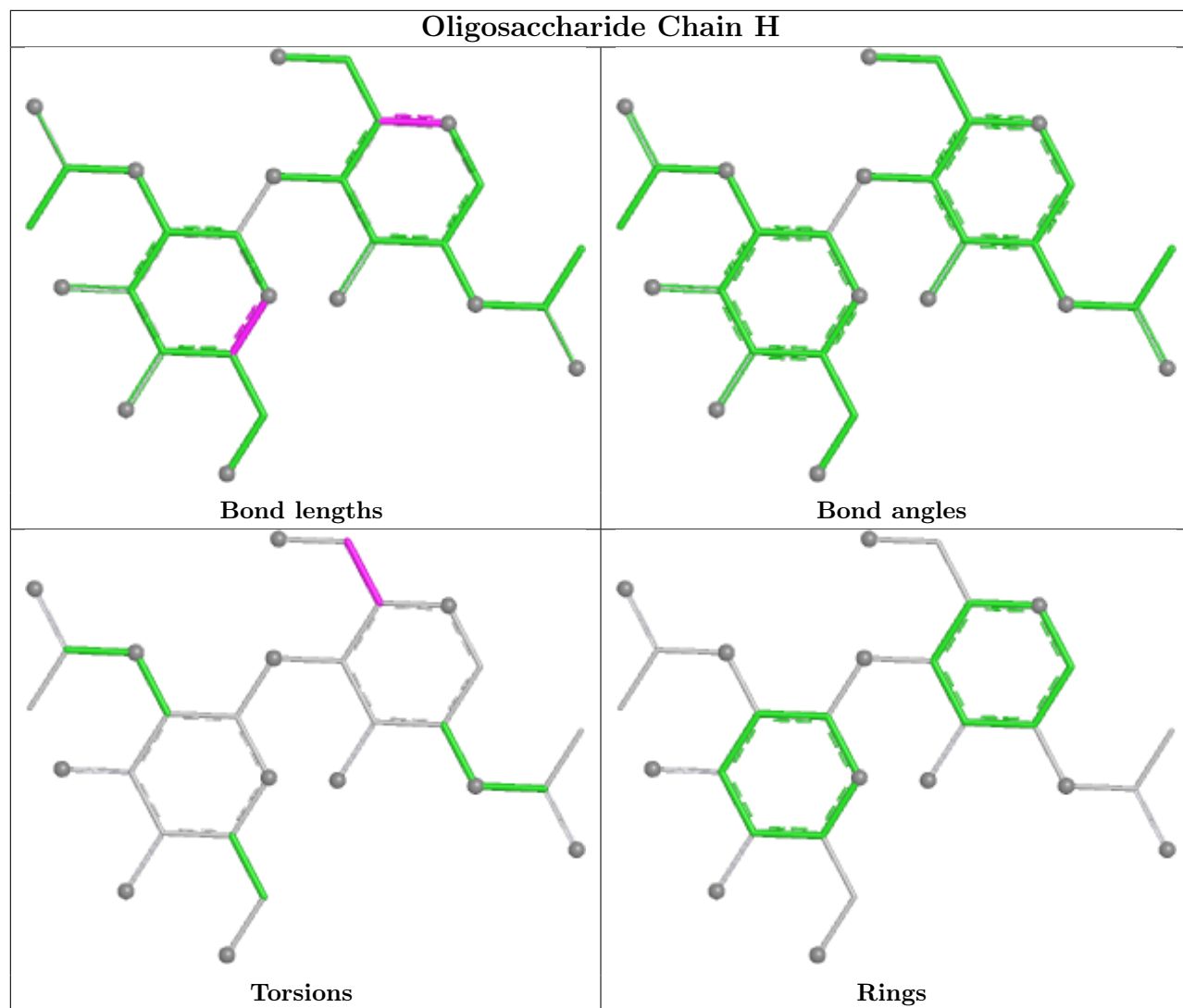
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

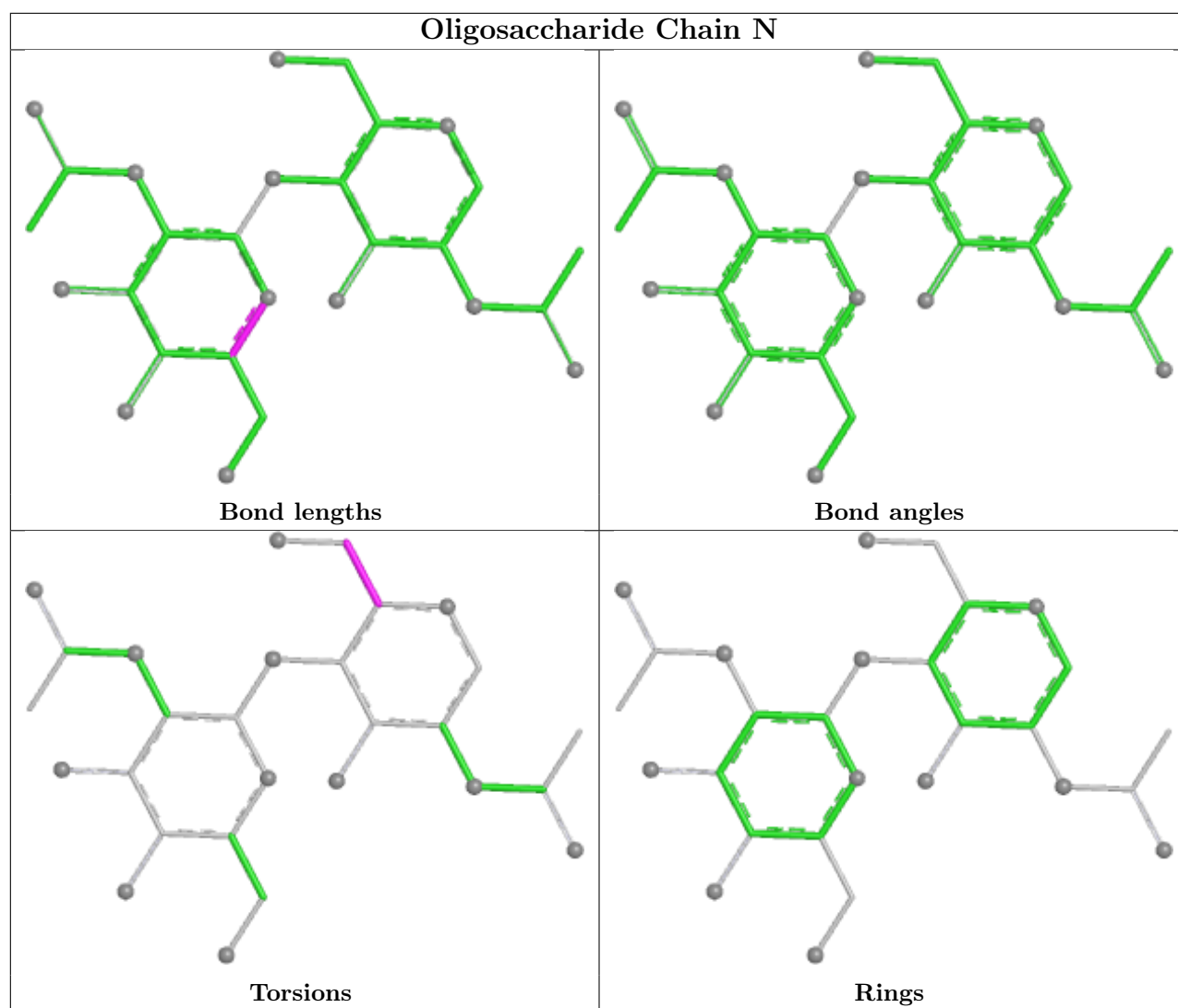


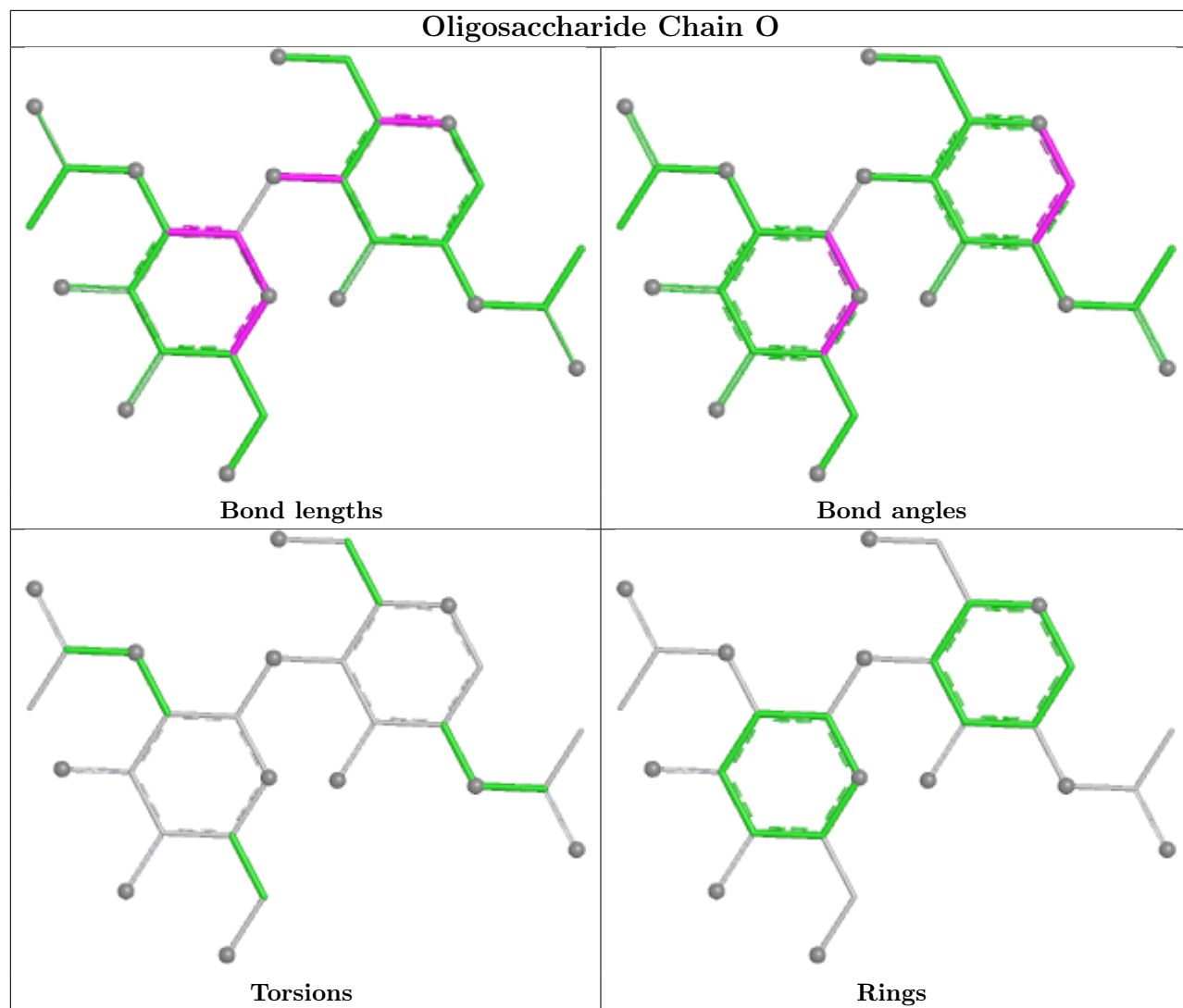


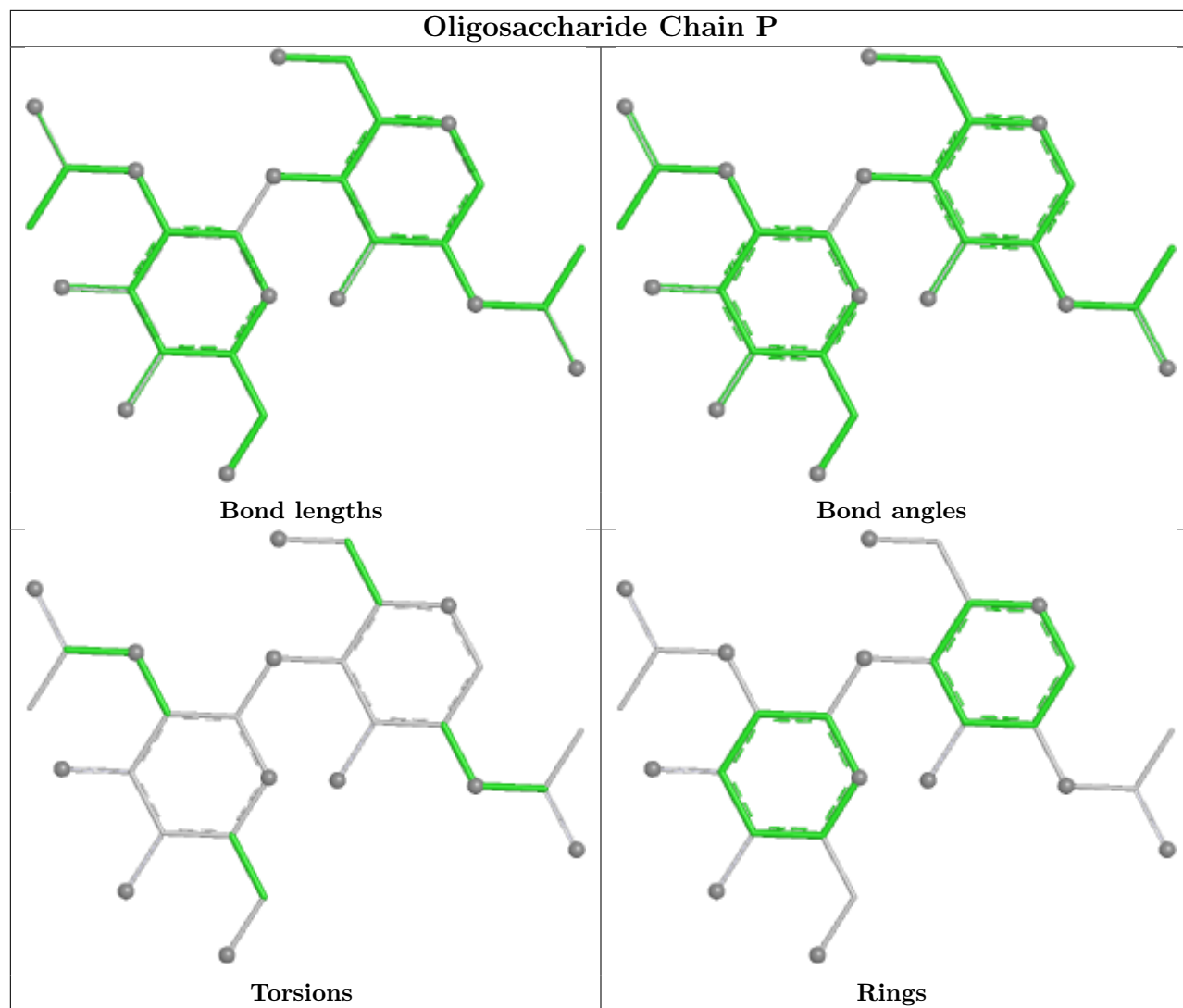


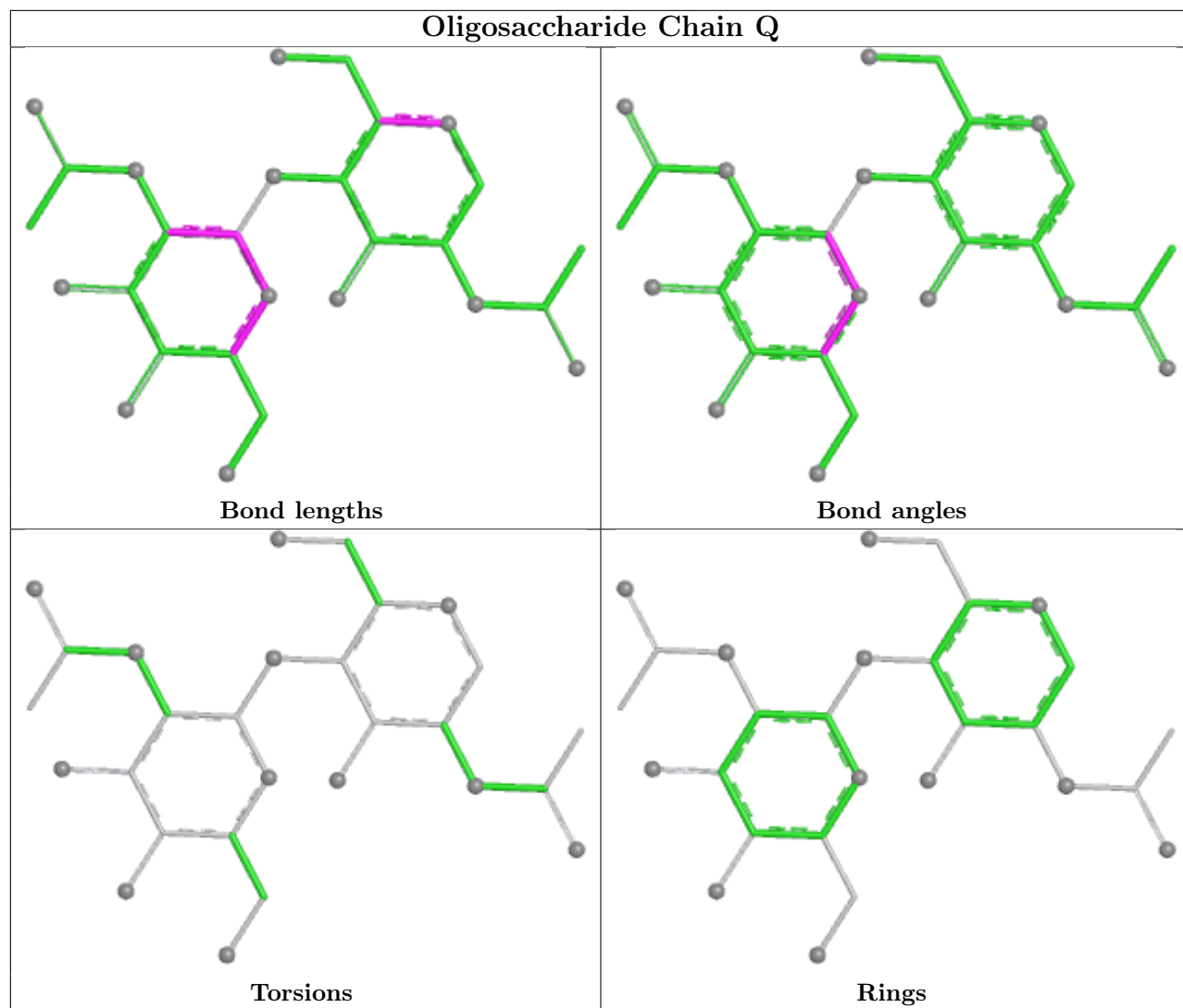


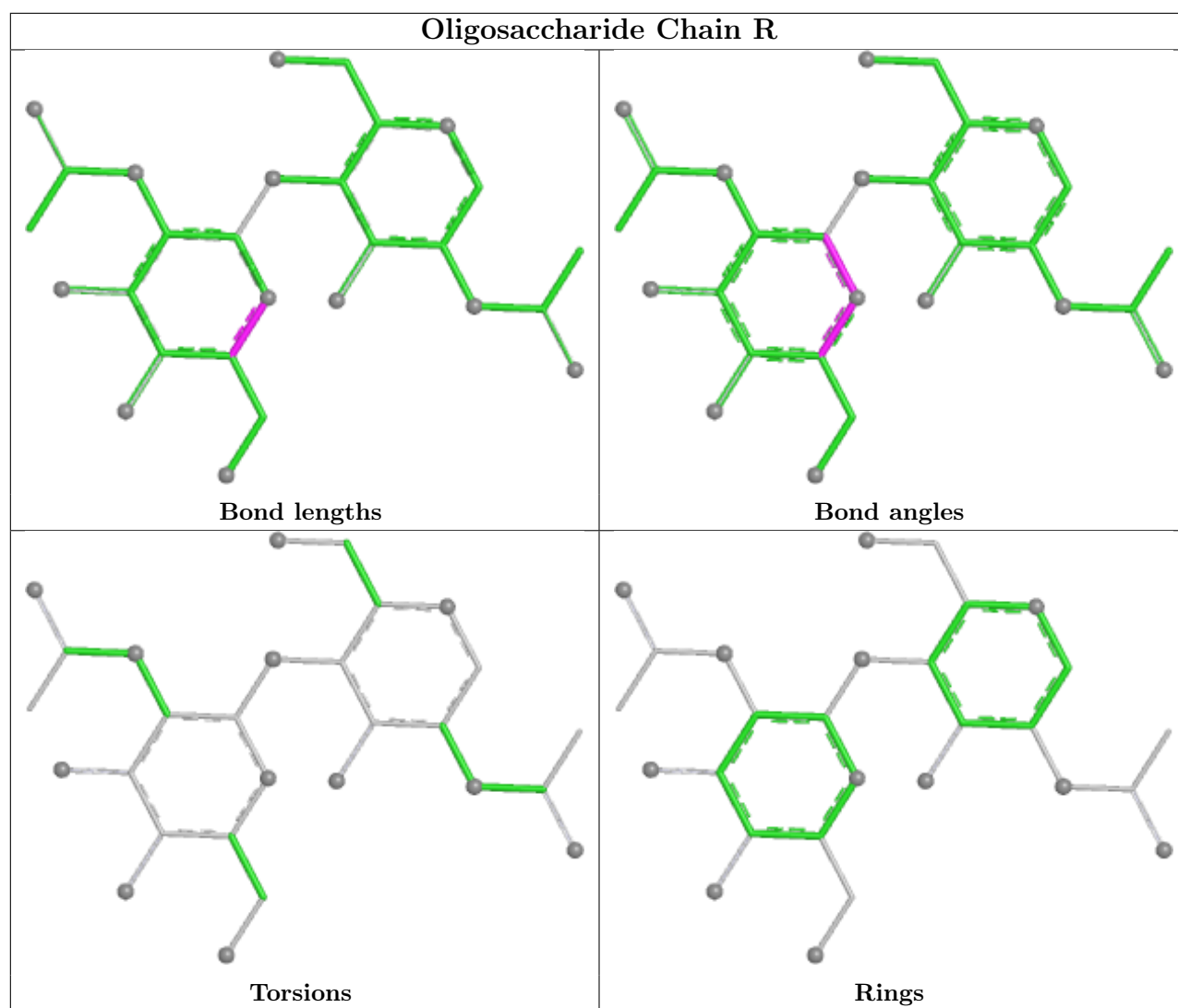


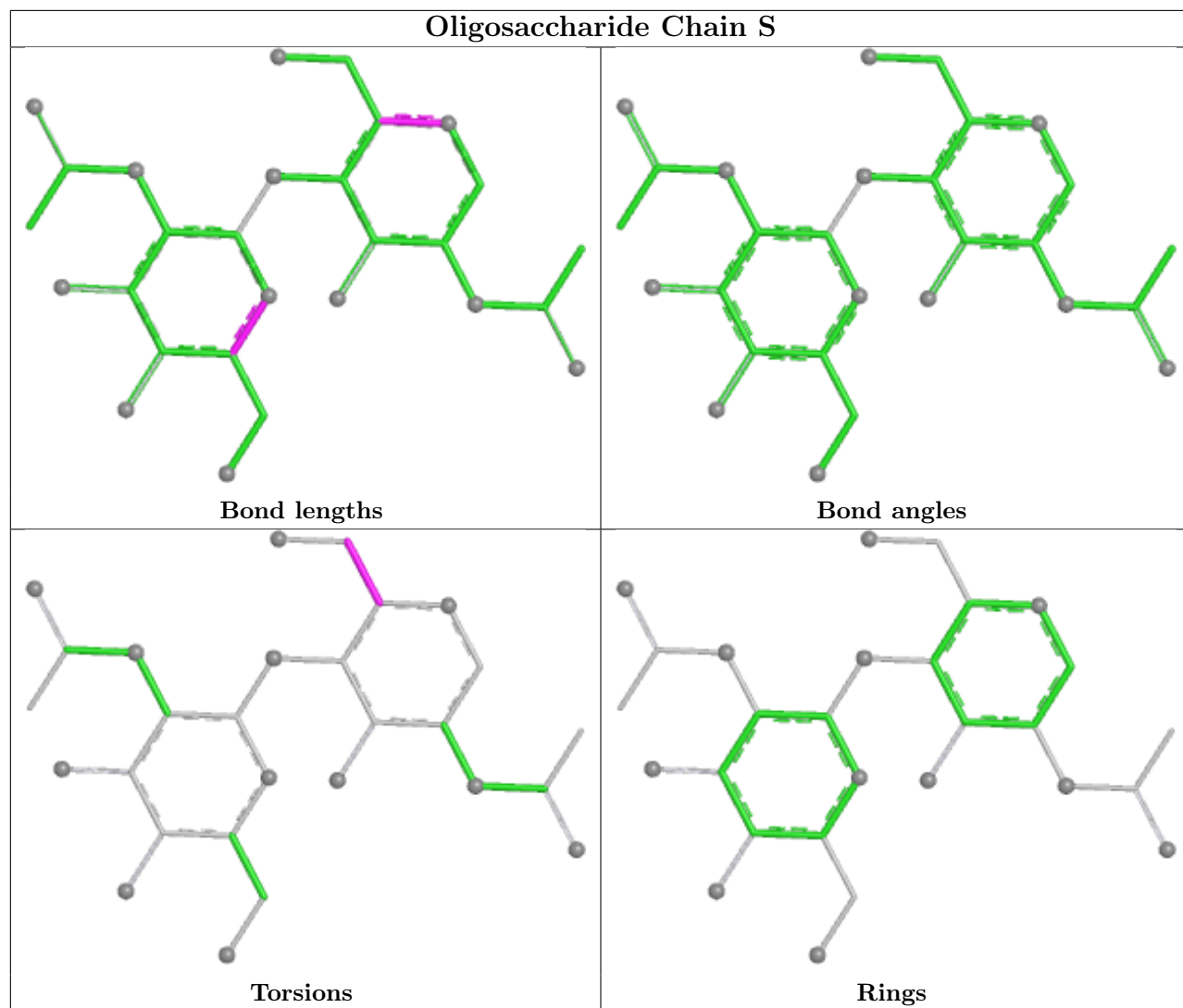


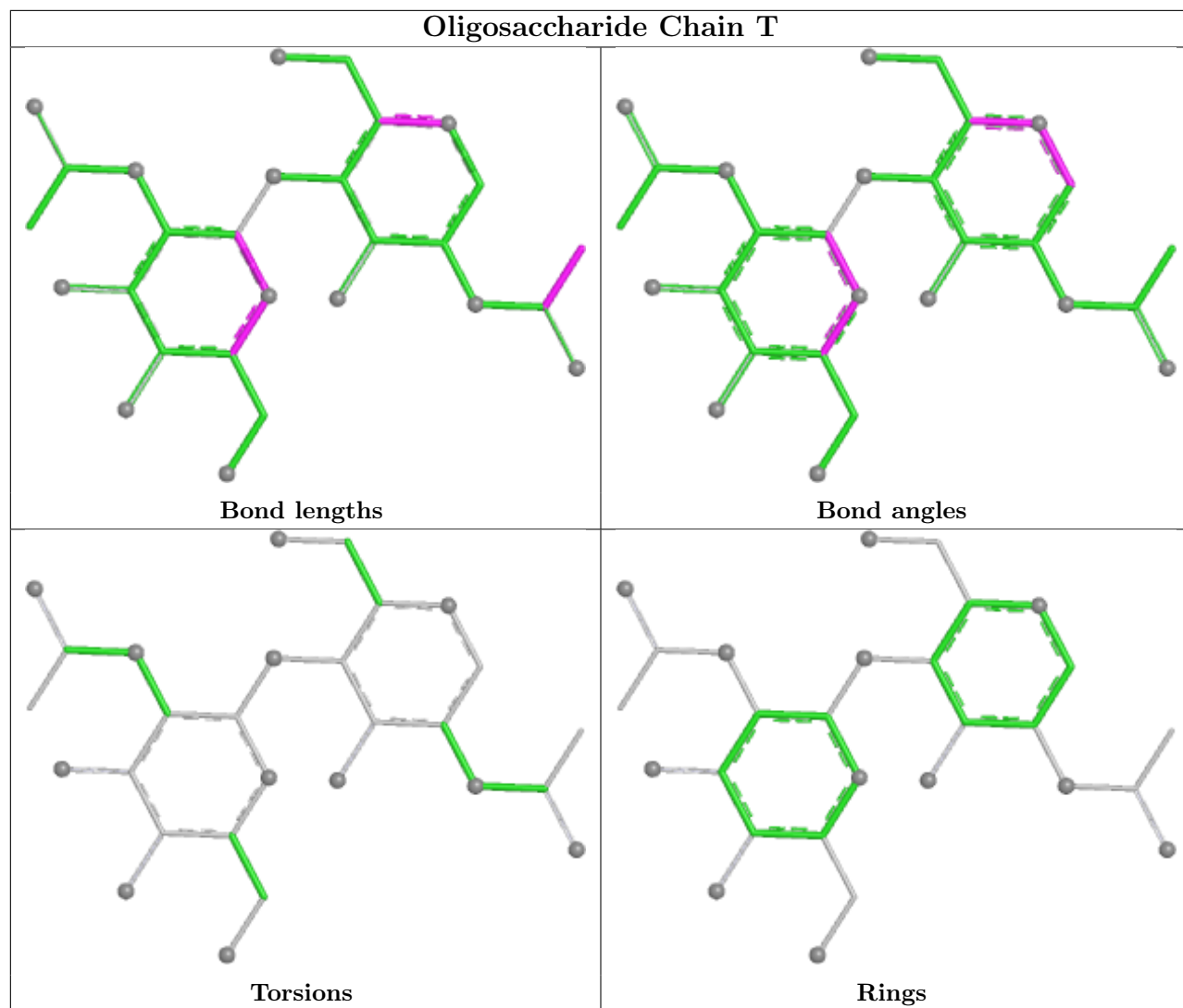


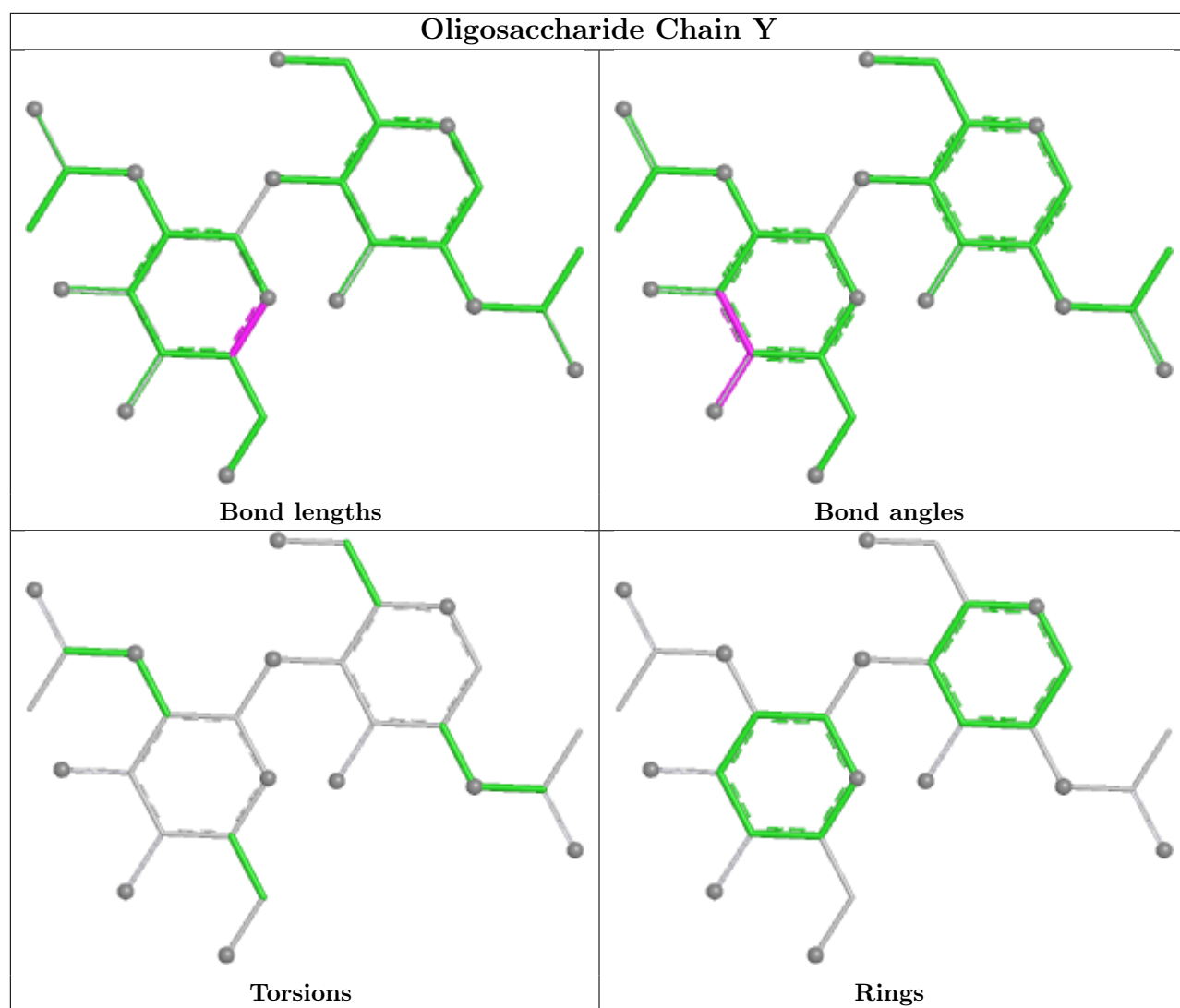


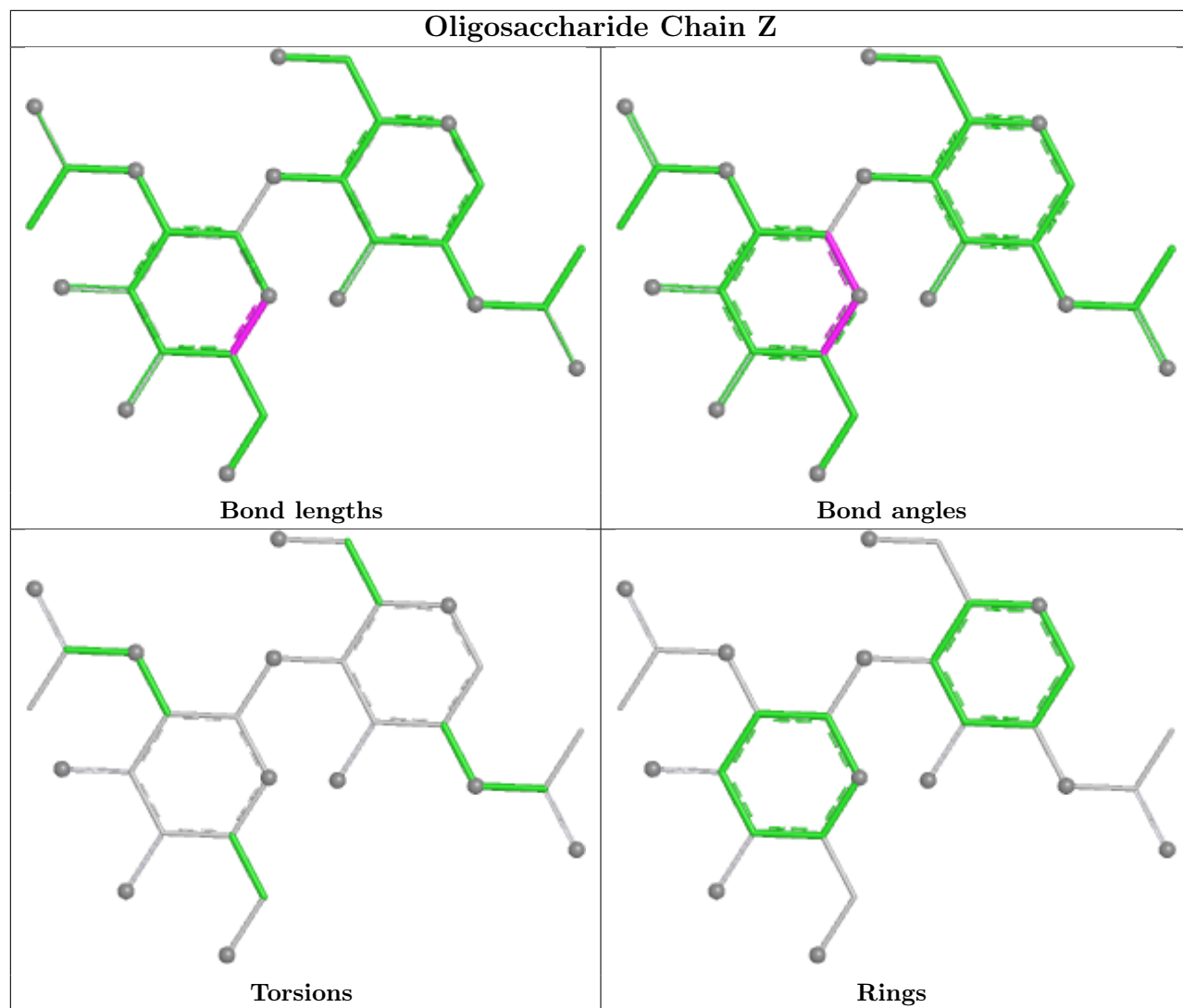


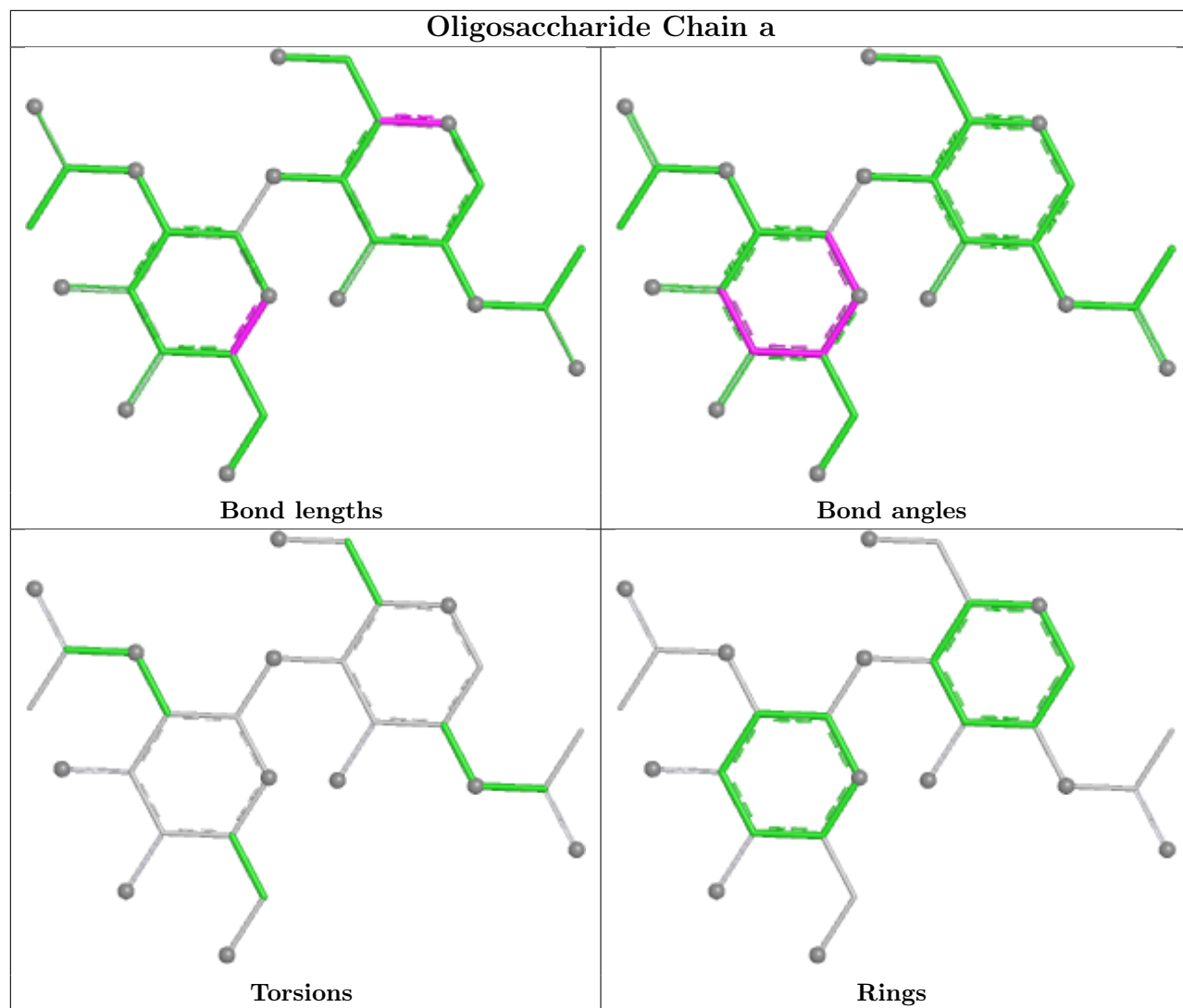


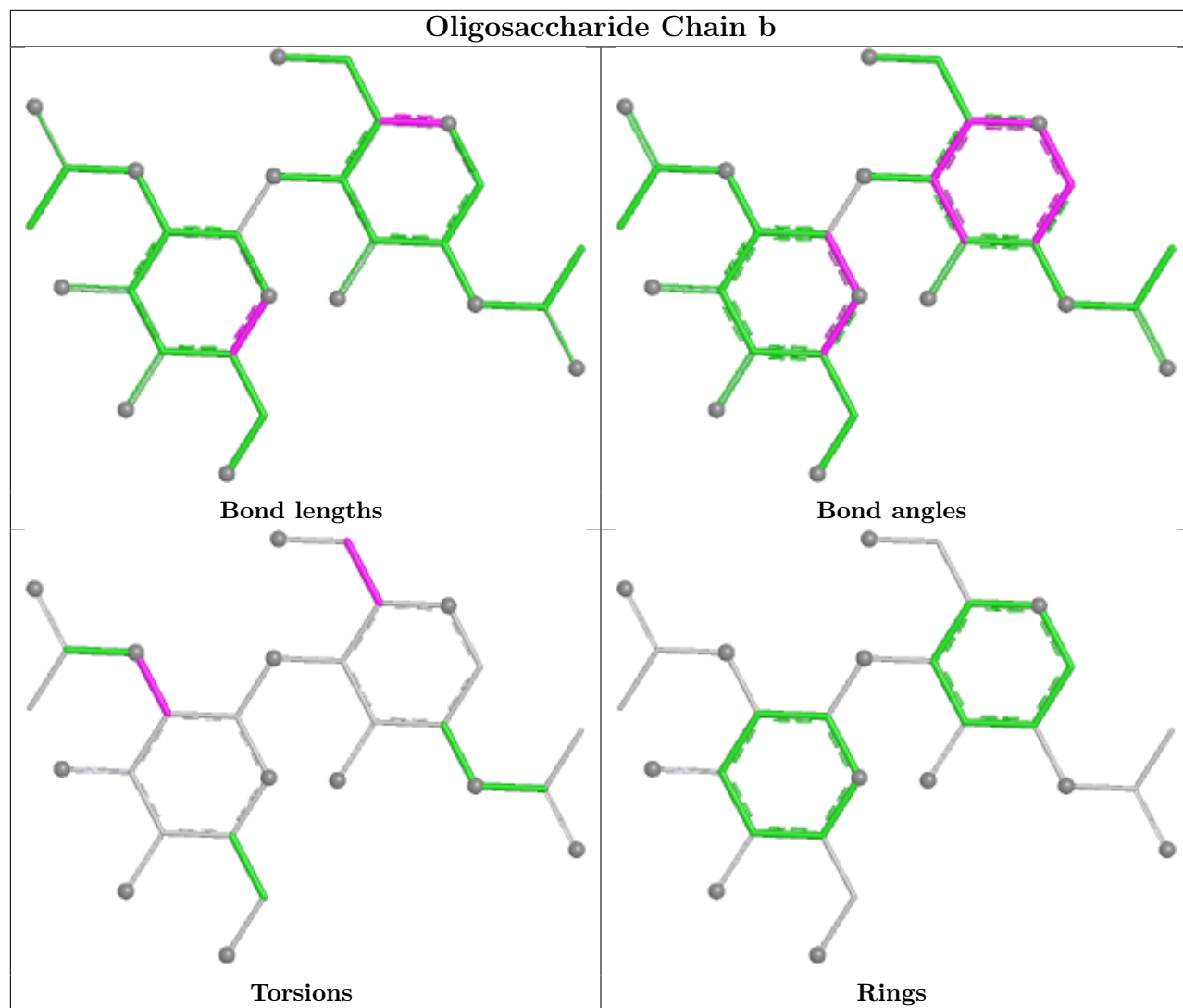


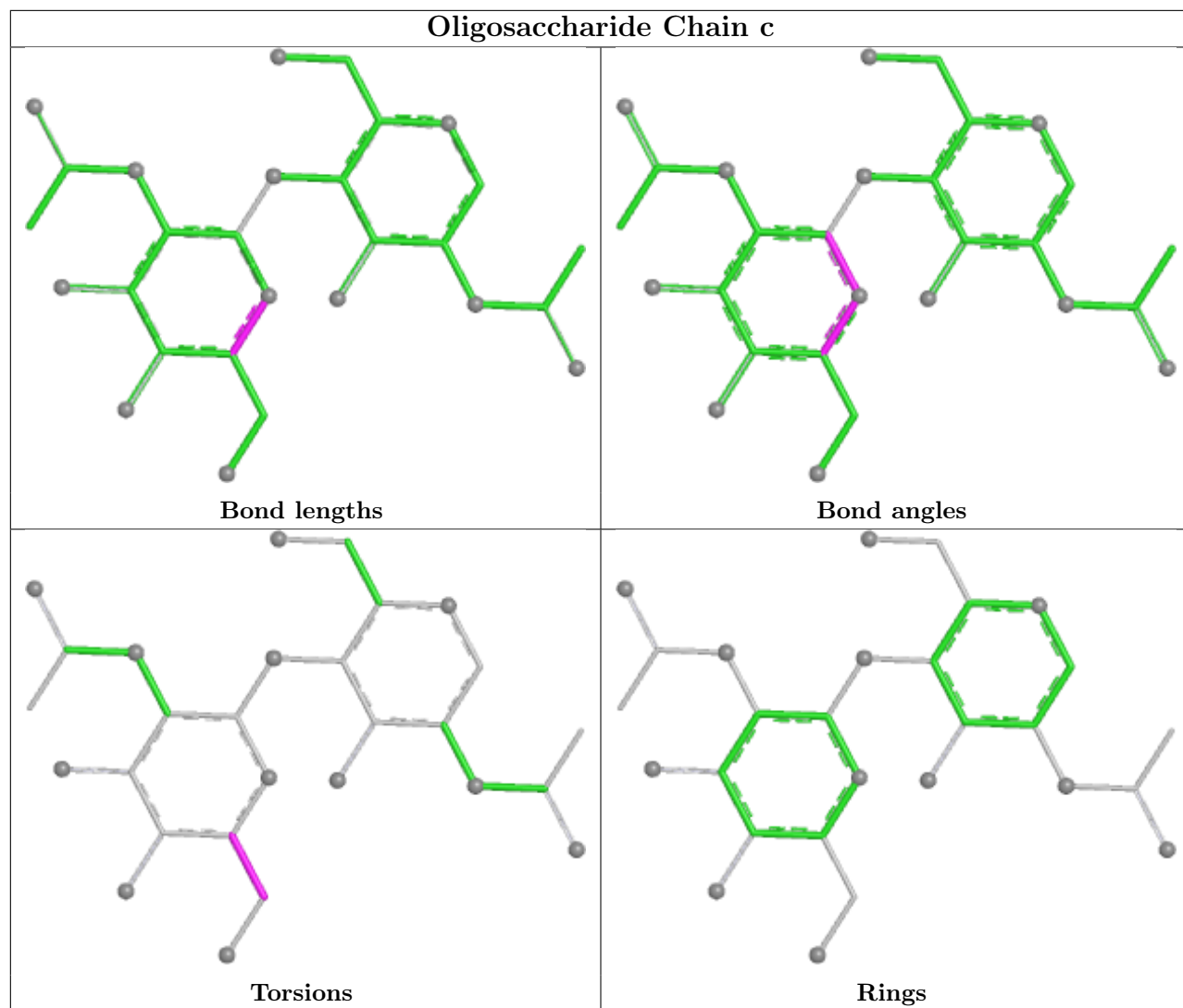


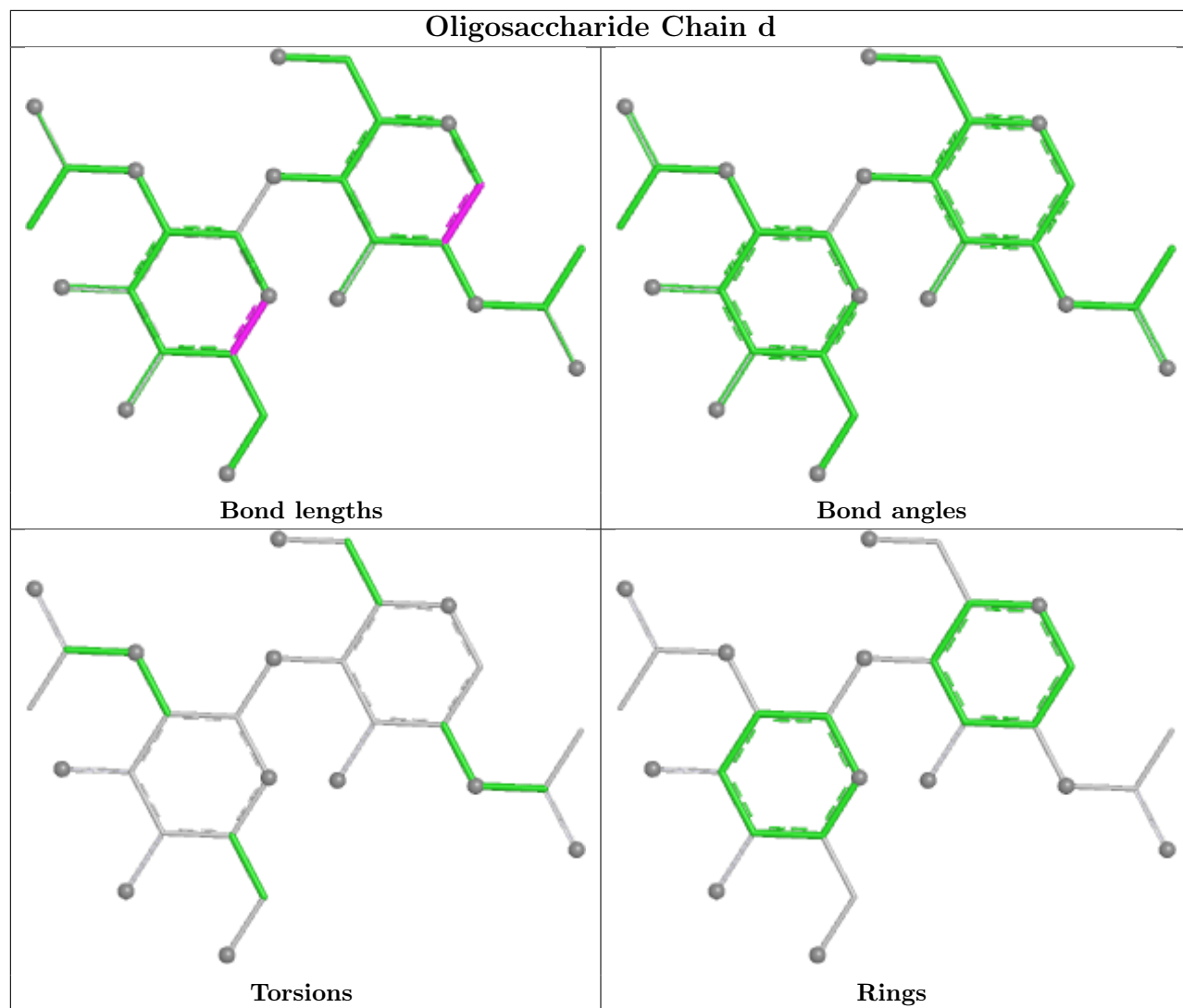


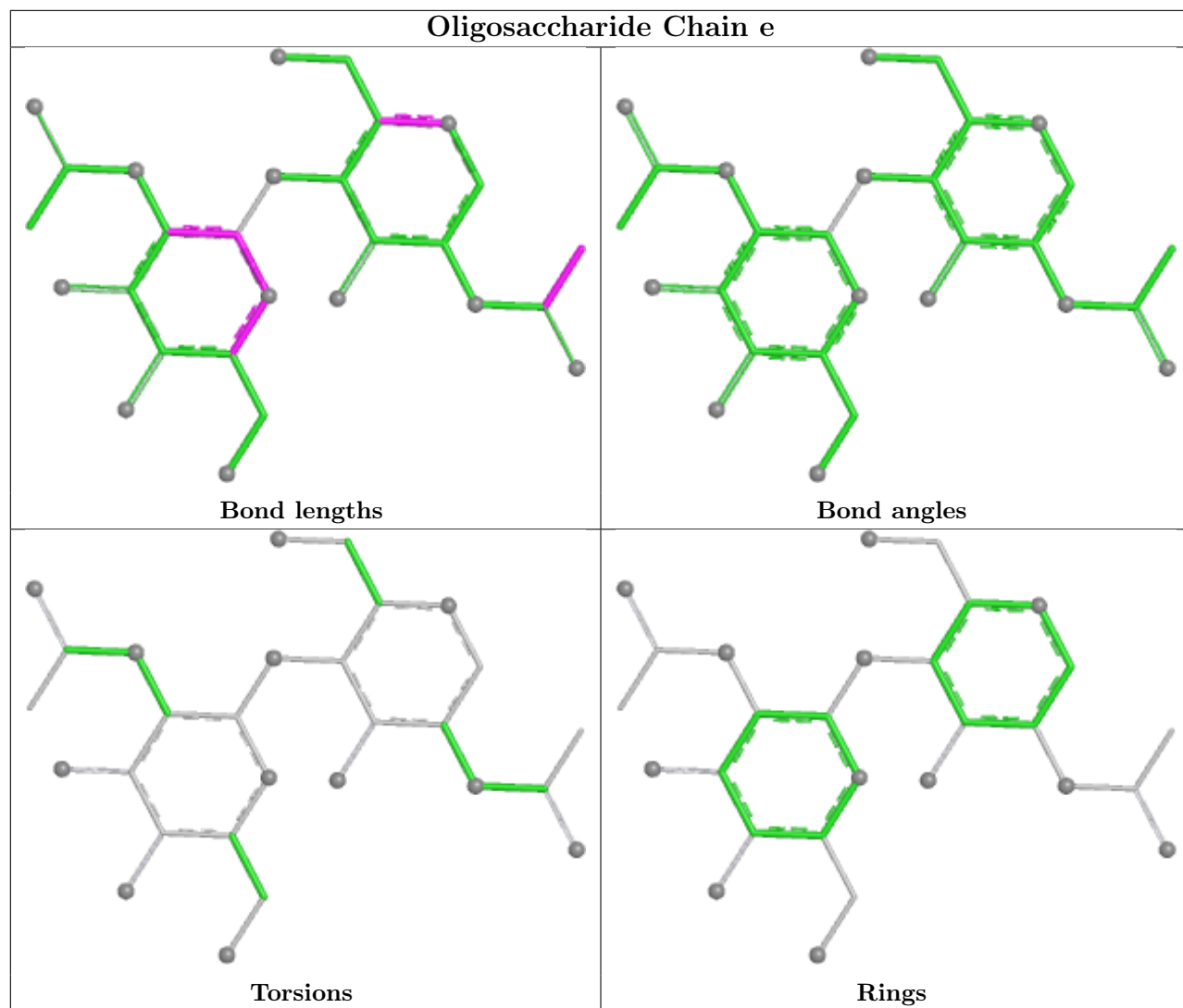


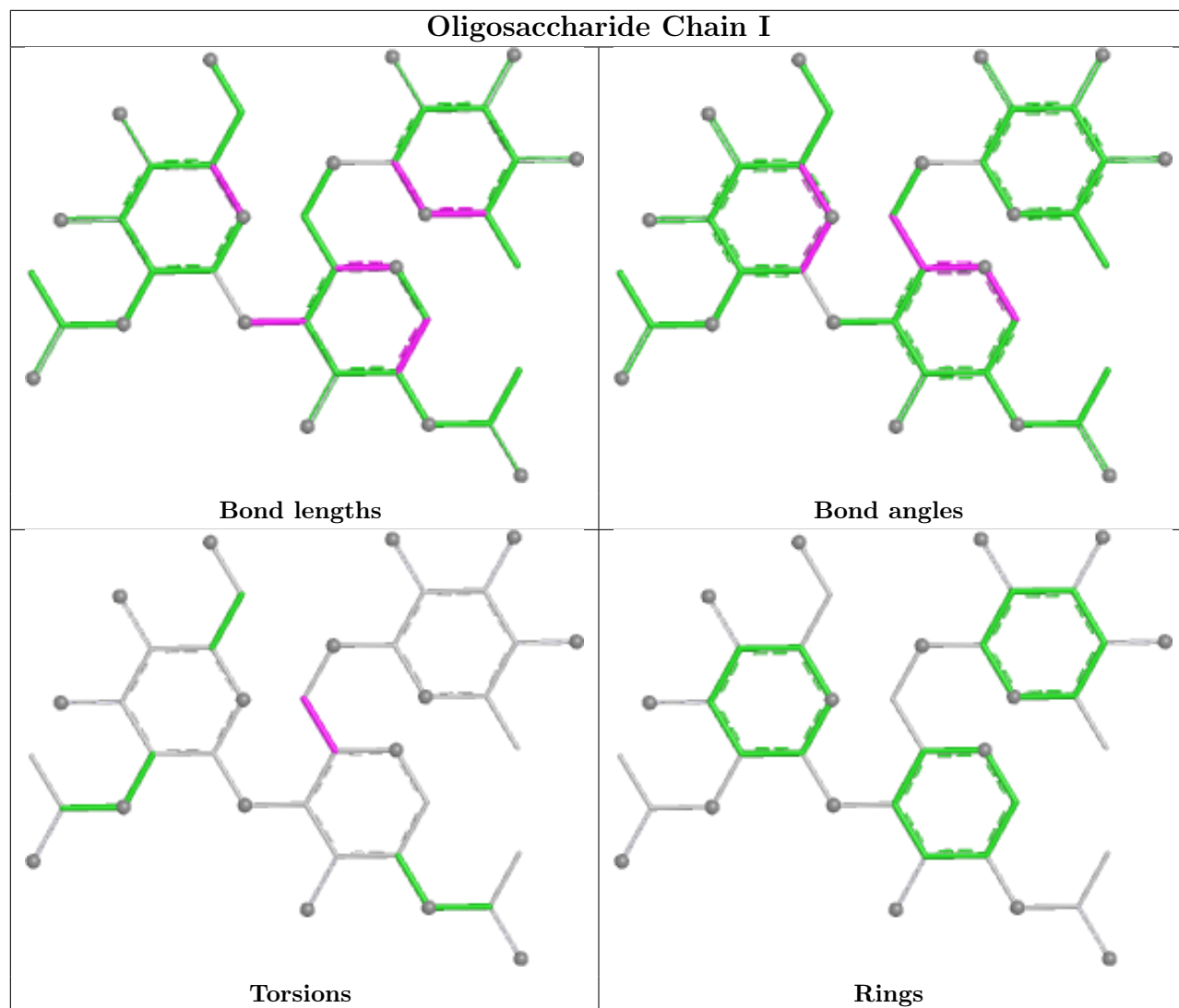


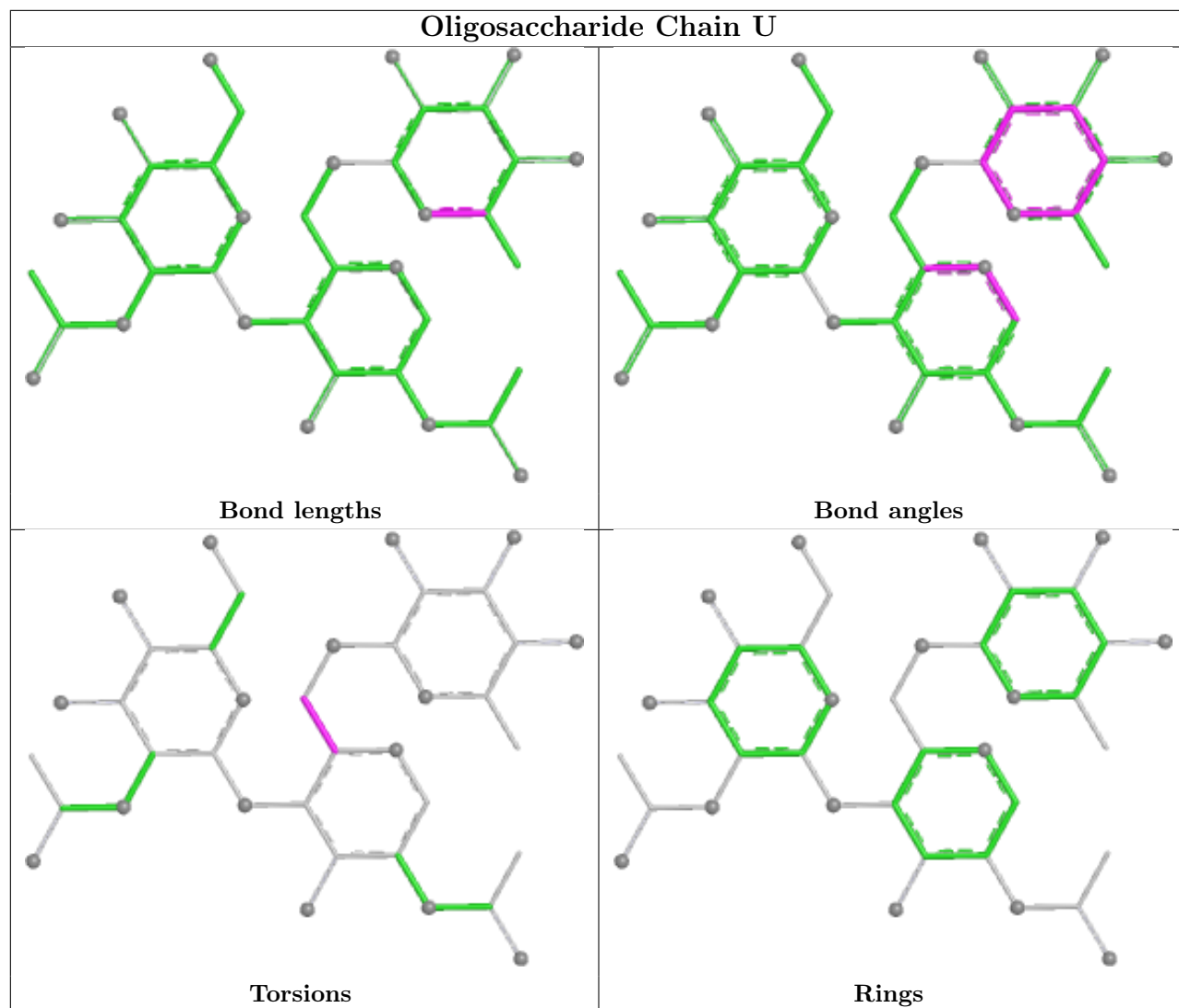


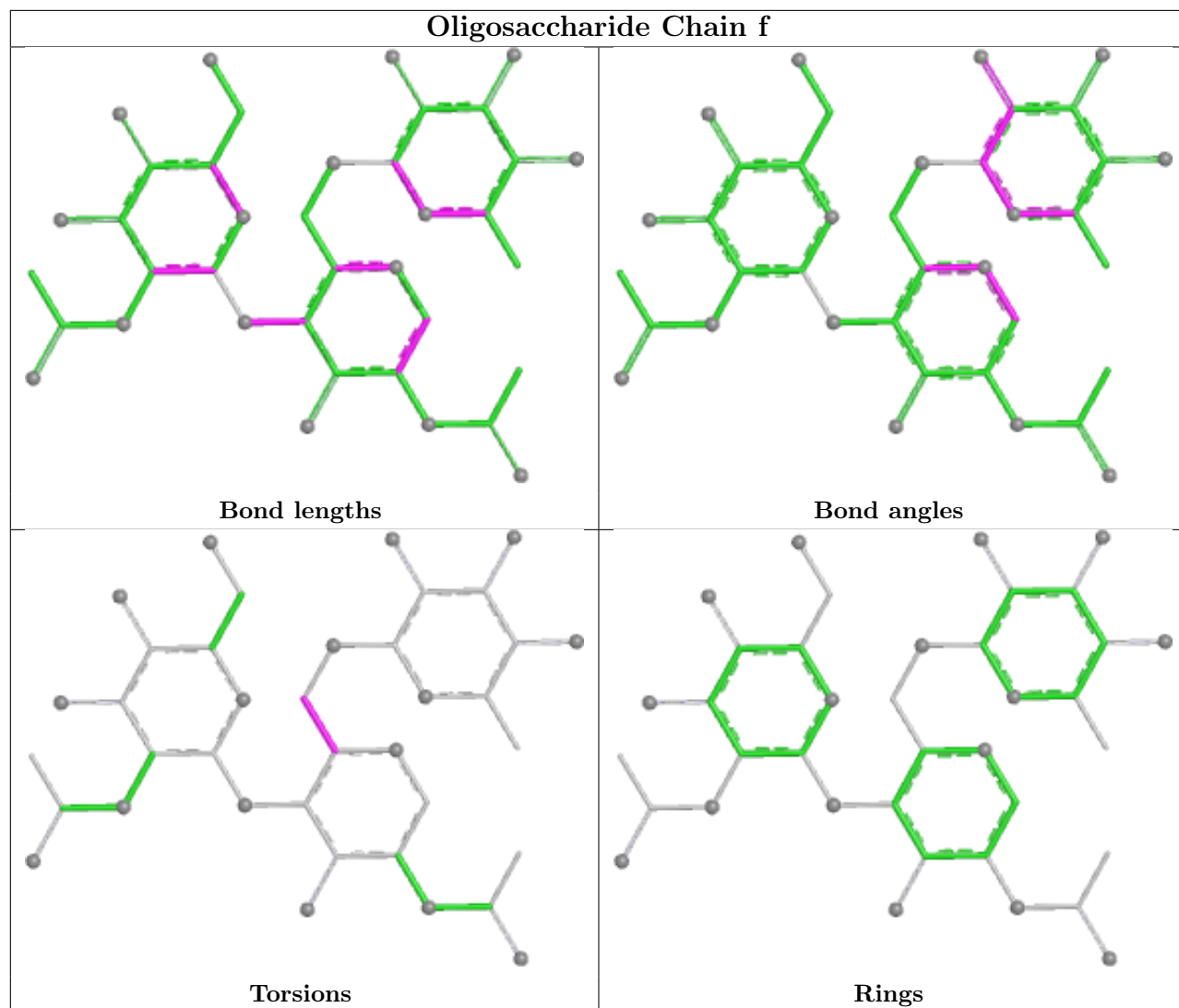


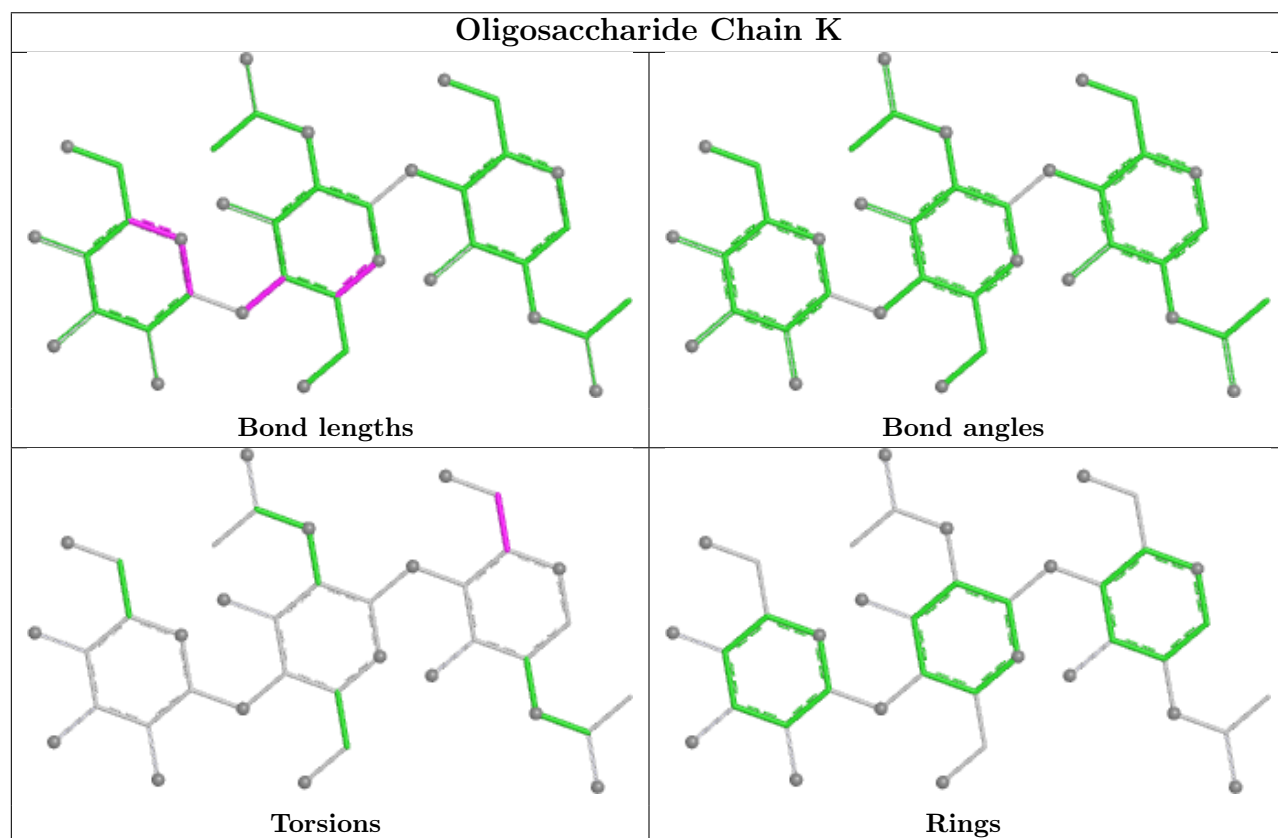
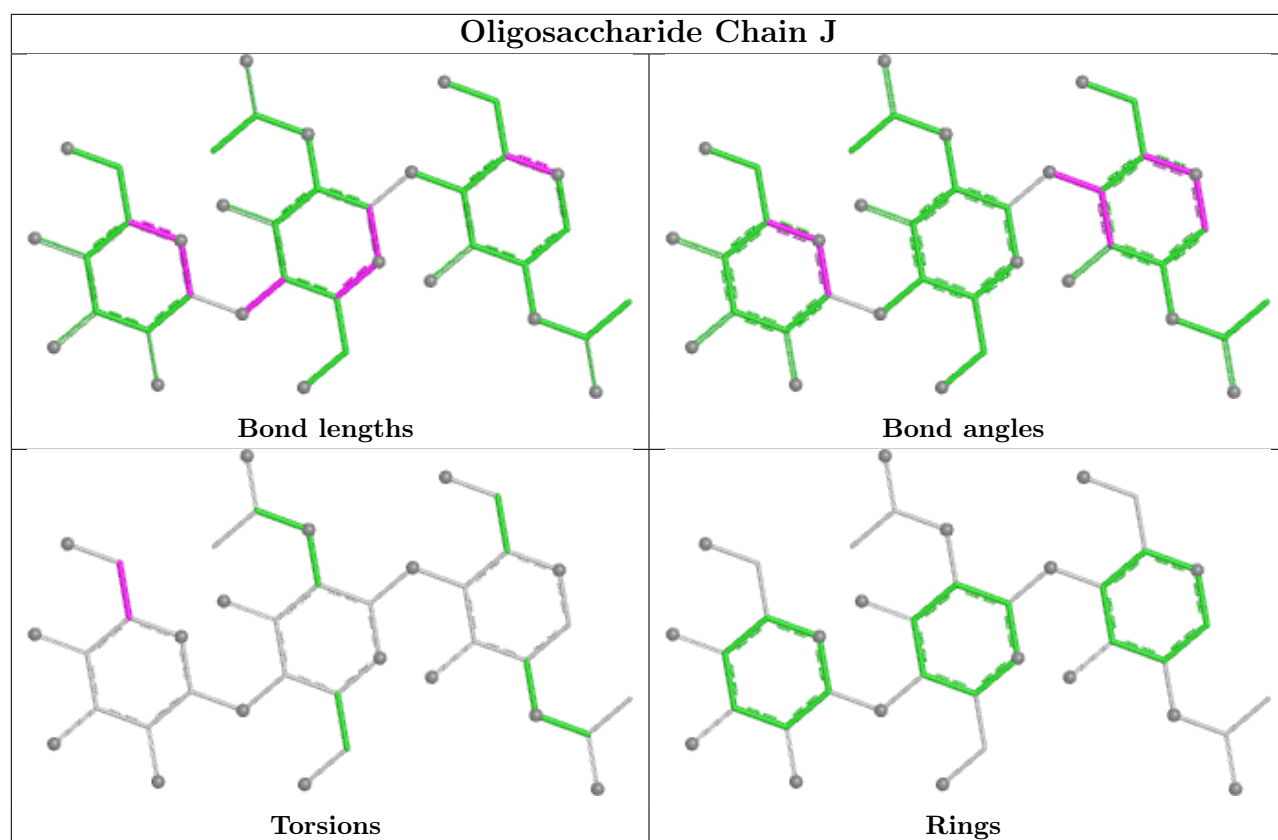


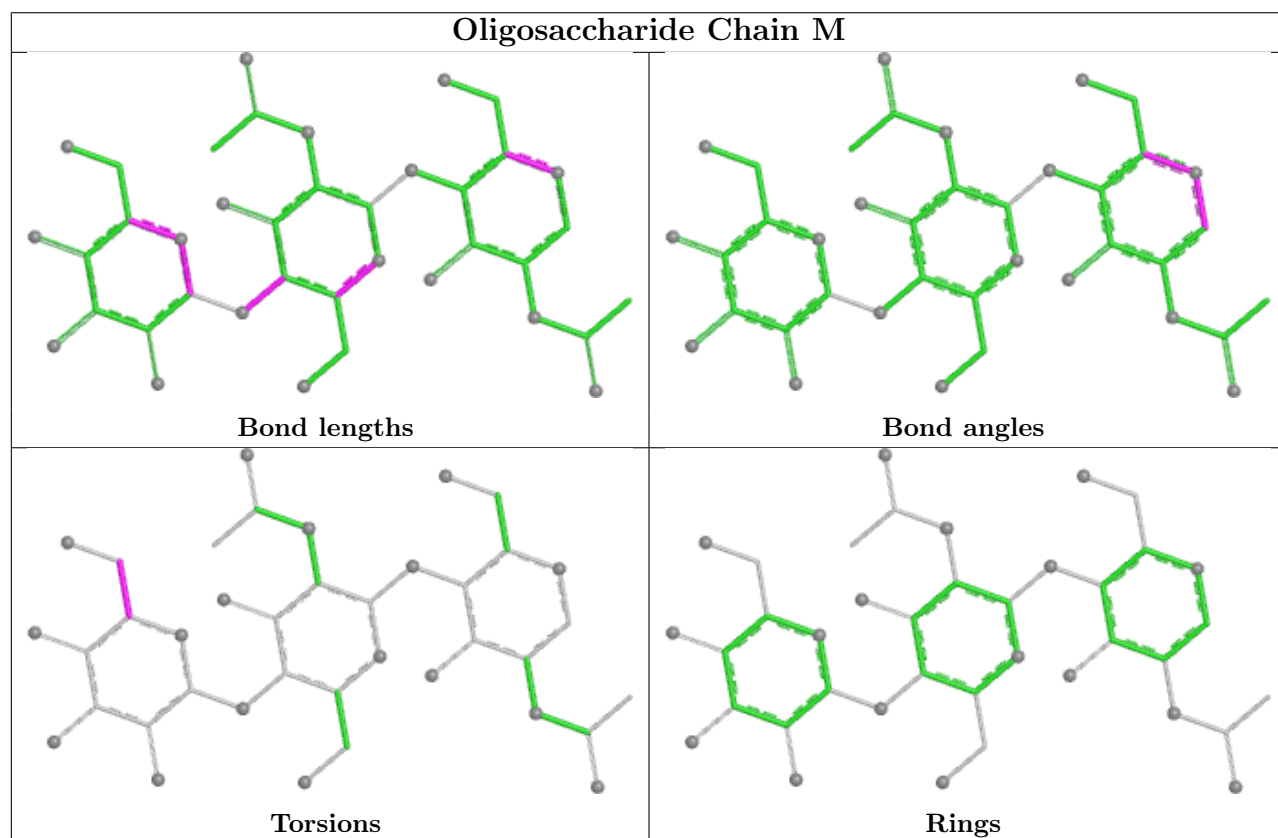
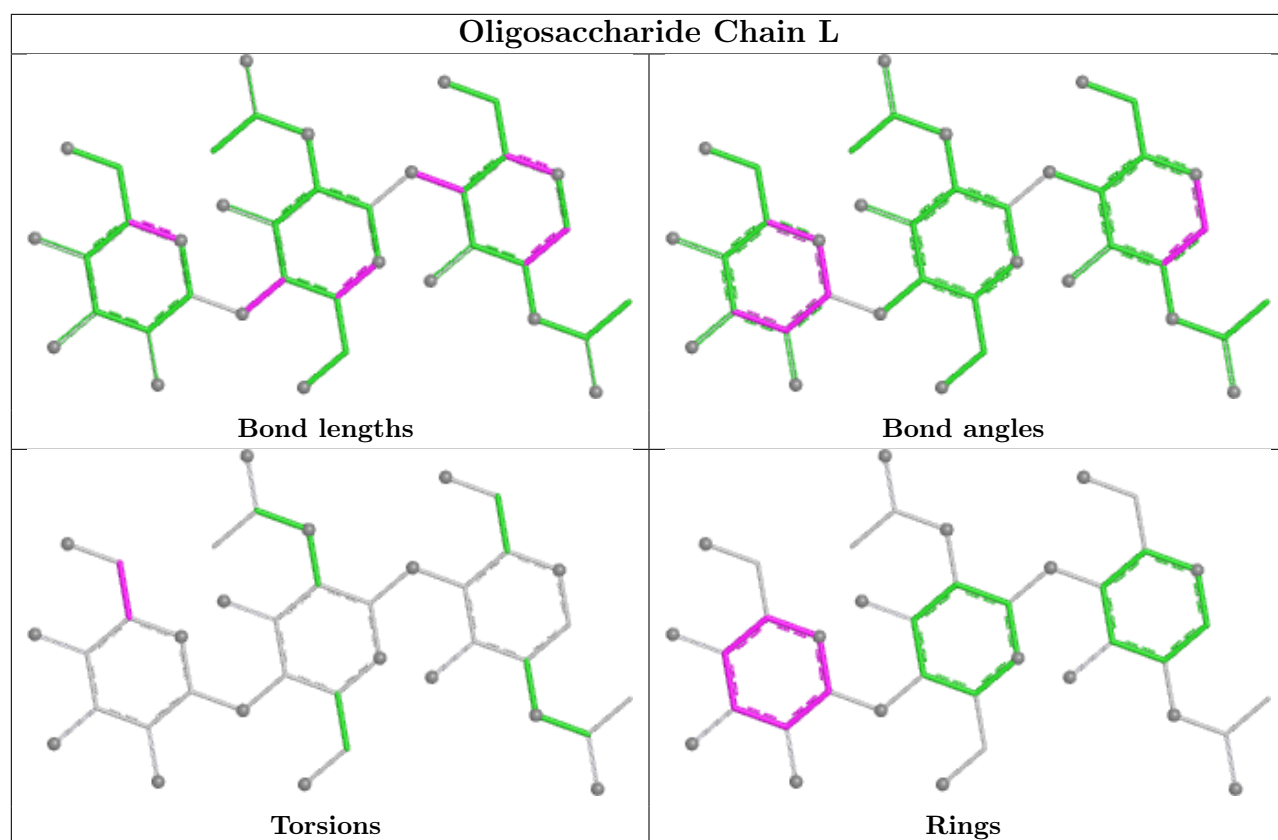


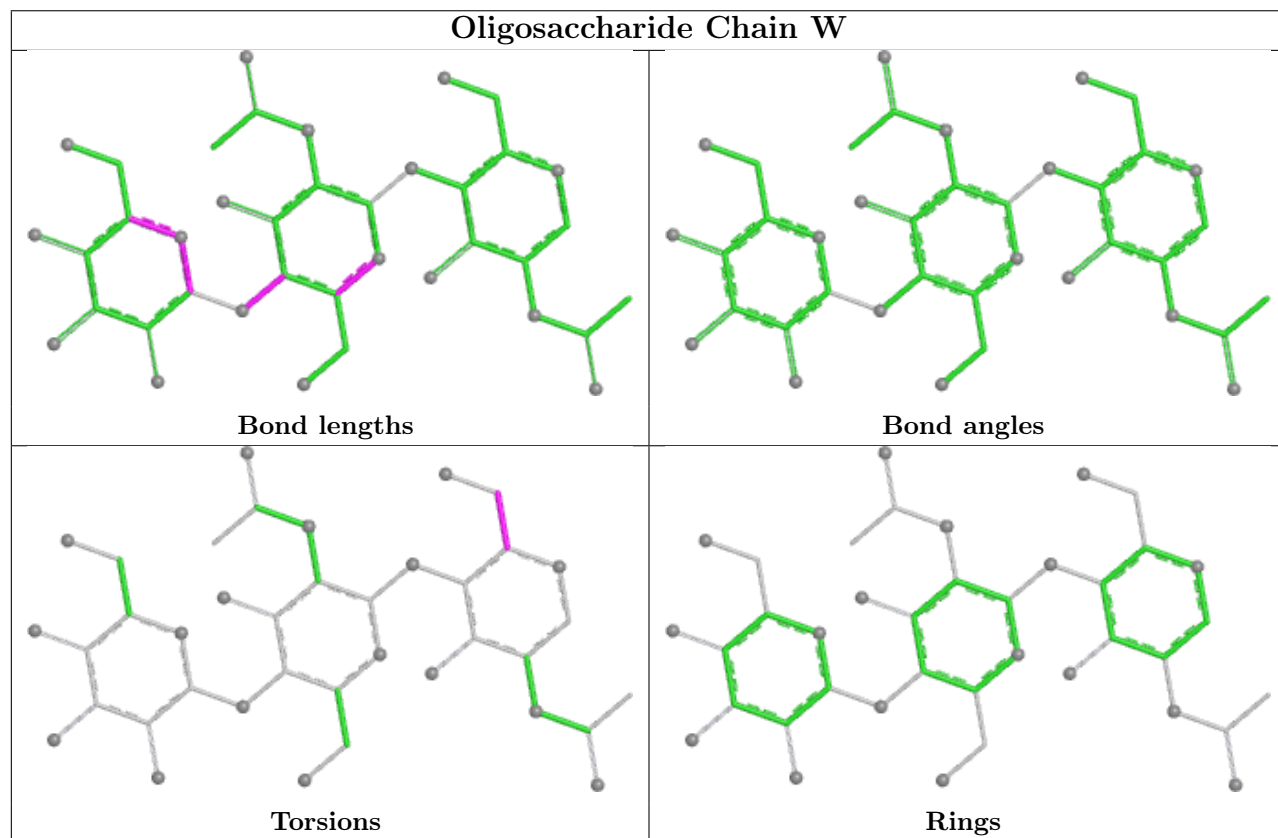
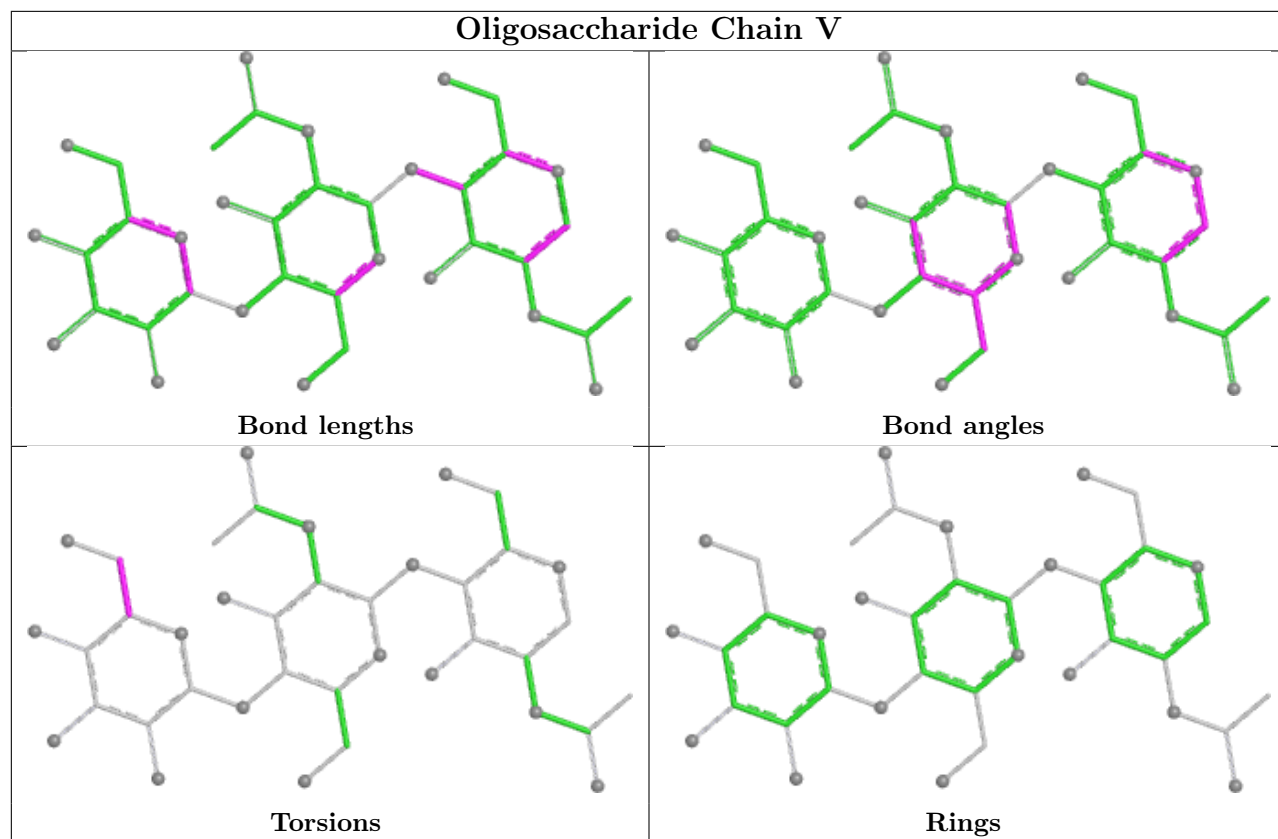


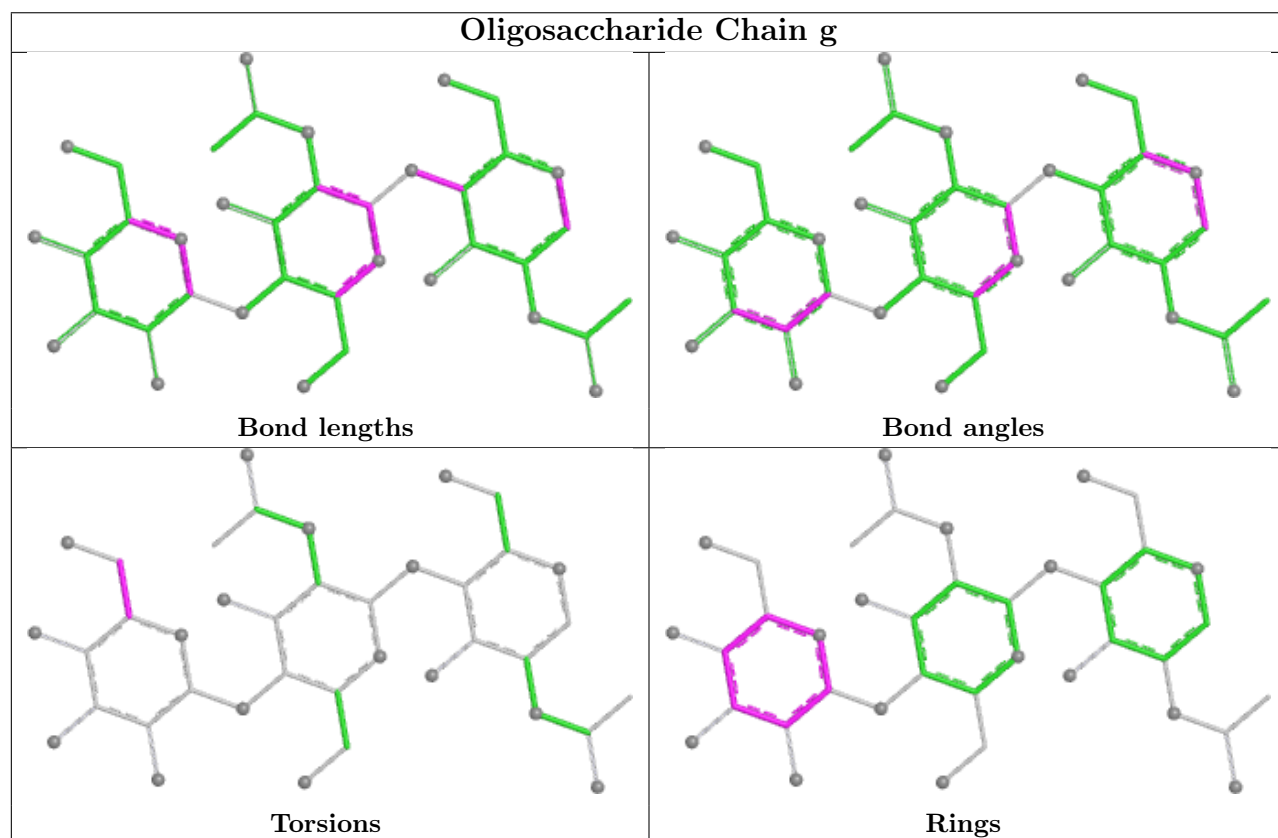
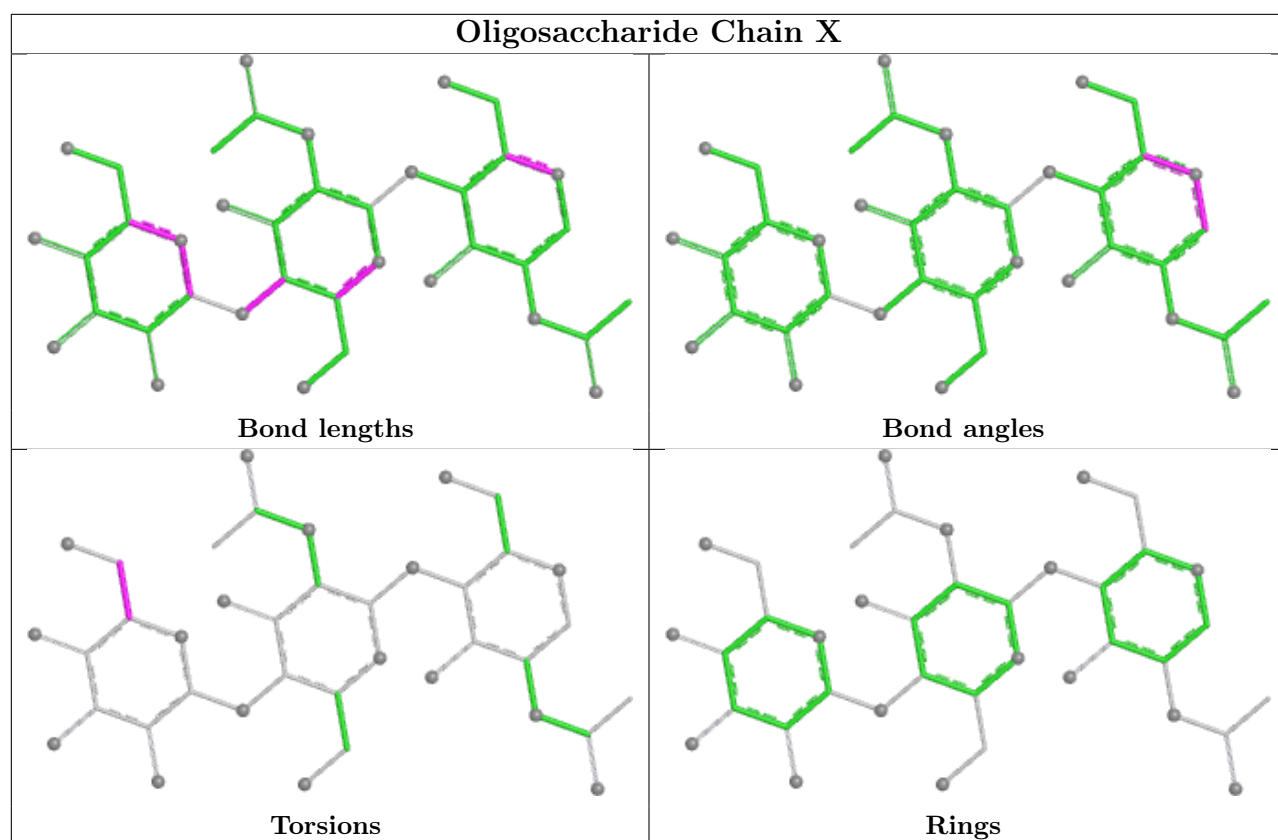


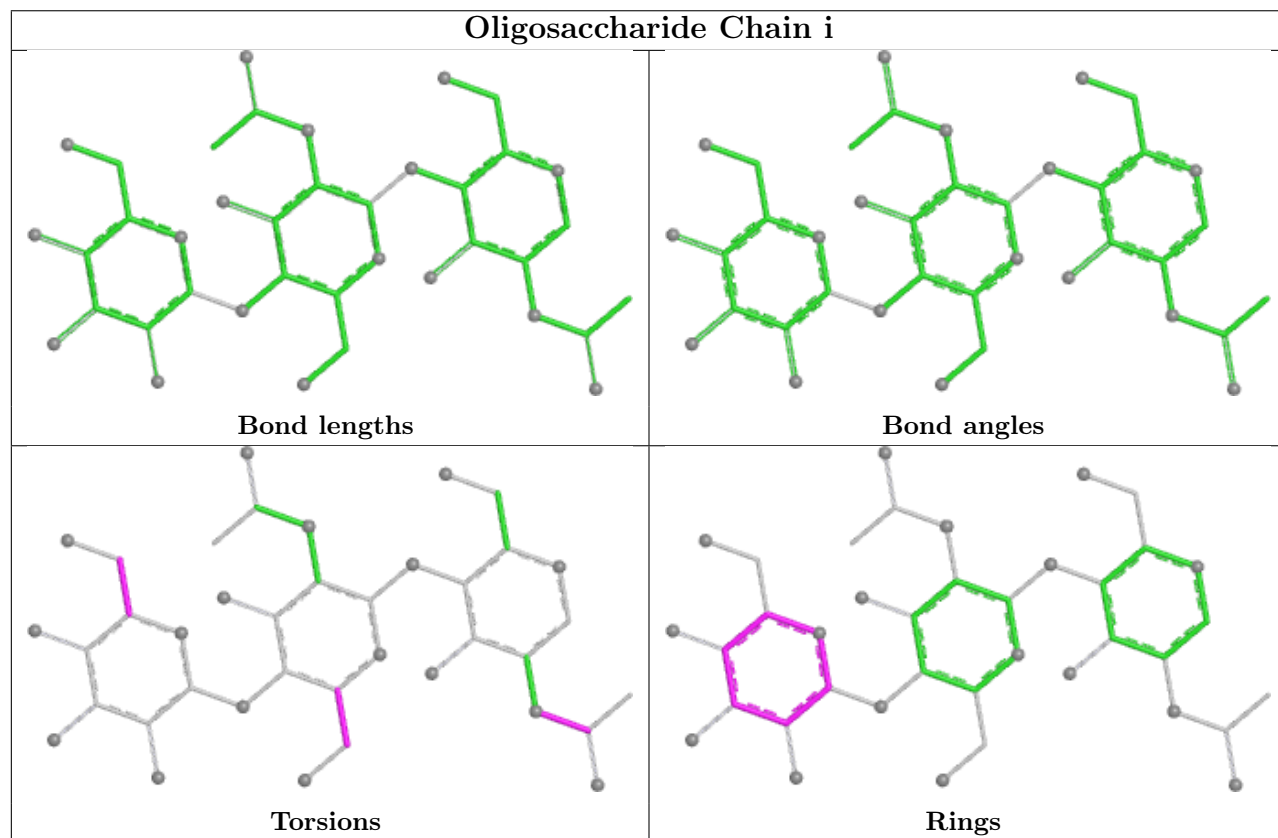
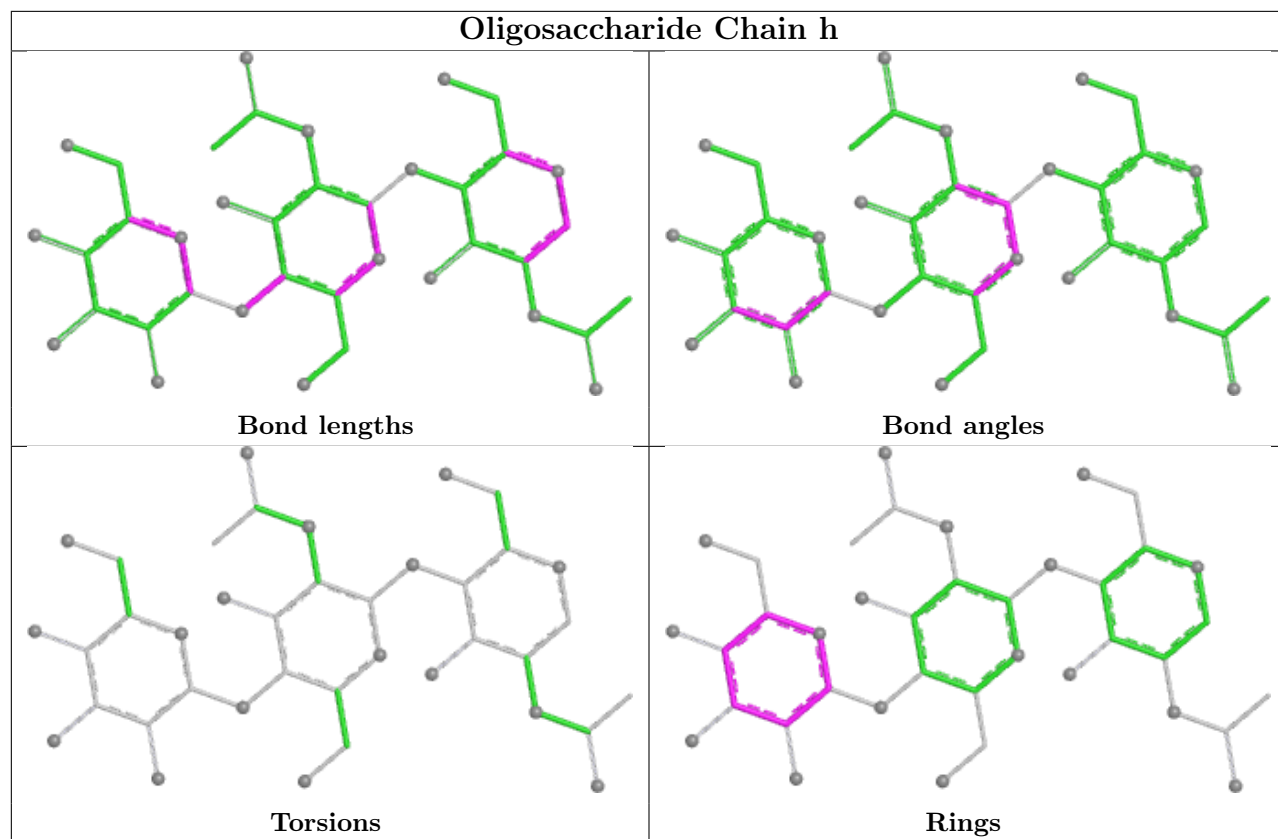


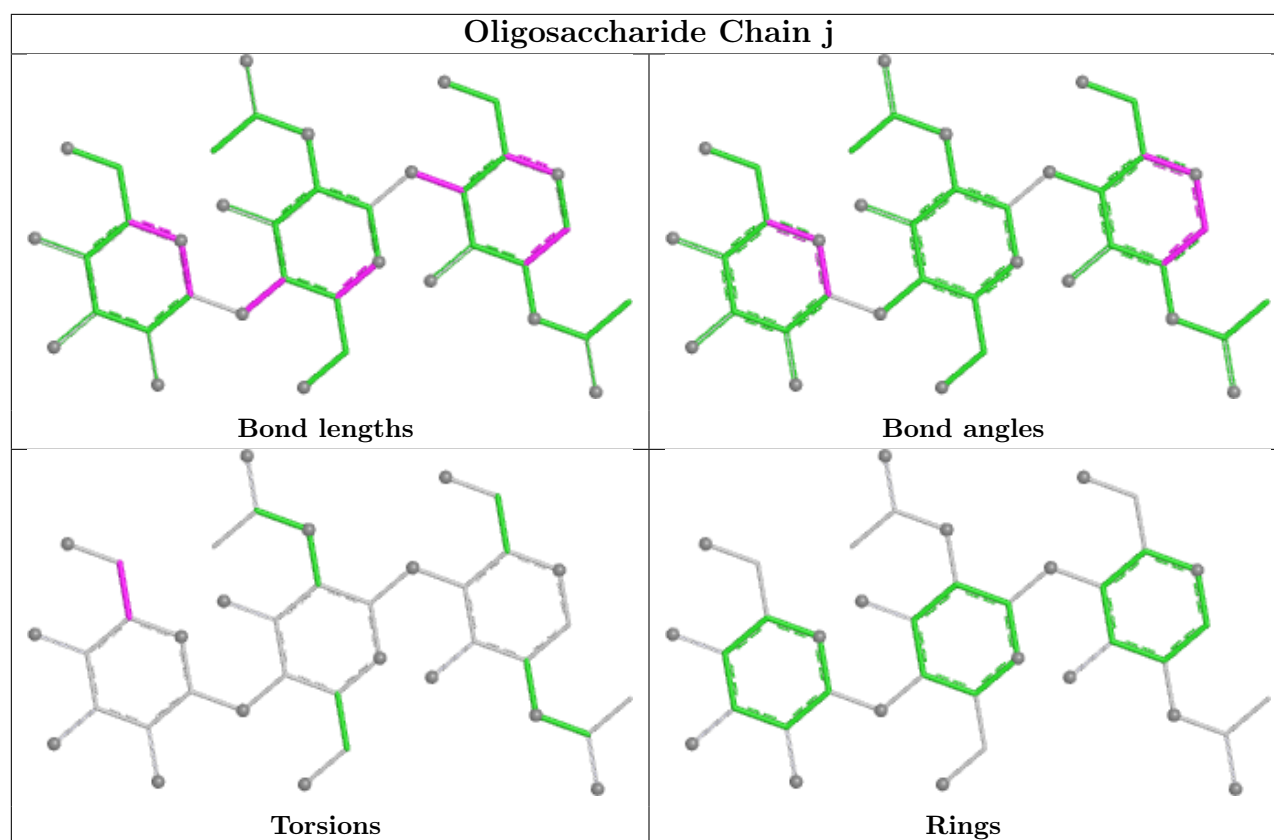












5.6 Ligand geometry [i](#)

24 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	$\# Z > 2$
5	NAG	A	1404	1	14,14,15	0.32	0	17,19,21	0.60	0
5	NAG	C	1404	1	14,14,15	0.23	0	17,19,21	0.44	0
5	NAG	A	1402	1	14,14,15	1.53	3 (21%)	17,19,21	1.16	1 (5%)
5	NAG	A	1409	1	14,14,15	1.34	2 (14%)	17,19,21	0.97	0
5	NAG	B	1403	1	14,14,15	1.12	1 (7%)	17,19,21	0.90	1 (5%)
5	NAG	B	1404	1	14,14,15	1.14	1 (7%)	17,19,21	1.10	1 (5%)
5	NAG	C	1406	1	14,14,15	1.49	3 (21%)	17,19,21	0.63	0
5	NAG	A	1401	1	14,14,15	1.42	2 (14%)	17,19,21	1.43	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	A	1405	1	14,14,15	0.34	0	17,19,21	0.45	0
5	NAG	A	1408	1	14,14,15	1.32	2 (14%)	17,19,21	0.91	1 (5%)
5	NAG	A	1406	1	14,14,15	1.31	2 (14%)	17,19,21	0.65	0
5	NAG	C	1407	1	14,14,15	1.49	2 (14%)	17,19,21	0.78	1 (5%)
5	NAG	A	1407	1	14,14,15	0.28	0	17,19,21	0.38	0
5	NAG	B	1406	1	14,14,15	0.28	0	17,19,21	0.37	0
5	NAG	A	1403	1	14,14,15	0.50	0	17,19,21	0.65	1 (5%)
5	NAG	C	1403	1	14,14,15	1.11	1 (7%)	17,19,21	1.10	1 (5%)
5	NAG	B	1402	1	14,14,15	0.50	0	17,19,21	0.65	1 (5%)
5	NAG	B	1405	1	14,14,15	1.61	2 (14%)	17,19,21	0.69	0
5	NAG	B	1407	1	14,14,15	1.27	1 (7%)	17,19,21	1.02	2 (11%)
5	NAG	C	1402	1	14,14,15	1.16	1 (7%)	17,19,21	1.19	1 (5%)
5	NAG	B	1408	1	14,14,15	1.68	3 (21%)	17,19,21	0.95	0
5	NAG	C	1405	1	14,14,15	0.25	0	17,19,21	0.38	0
5	NAG	C	1401	1	14,14,15	0.64	0	17,19,21	0.43	0
5	NAG	B	1401	1	14,14,15	0.51	0	17,19,21	0.45	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
5	NAG	A	1404	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1404	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1402	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1409	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1403	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1404	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1406	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1401	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1405	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1408	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1406	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1407	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1407	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1406	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1403	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1403	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	B	1402	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1405	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1407	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1402	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1408	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1405	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1401	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1401	1	-	1/6/23/26	0/1/1/1

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1405	NAG	C1-C2	4.32	1.58	1.52
5	B	1404	NAG	O5-C1	-4.20	1.36	1.43
5	C	1403	NAG	O5-C1	-4.08	1.36	1.43
5	B	1408	NAG	O5-C5	3.47	1.50	1.43
5	A	1402	NAG	C1-C2	3.24	1.56	1.52
5	B	1408	NAG	O5-C1	3.20	1.49	1.43
5	C	1407	NAG	O5-C5	3.14	1.49	1.43
5	B	1407	NAG	O5-C5	3.09	1.49	1.43
5	A	1401	NAG	C1-C2	3.06	1.56	1.52
5	C	1407	NAG	C1-C2	3.03	1.56	1.52
5	C	1406	NAG	C8-C7	2.92	1.56	1.50
5	C	1406	NAG	O5-C5	2.90	1.49	1.43
5	C	1402	NAG	O5-C5	2.70	1.48	1.43
5	B	1405	NAG	O5-C5	2.69	1.48	1.43
5	A	1408	NAG	C1-C2	2.67	1.56	1.52
5	A	1406	NAG	O5-C5	2.64	1.48	1.43
5	A	1409	NAG	O5-C5	2.63	1.48	1.43
5	A	1401	NAG	O5-C5	2.59	1.48	1.43
5	A	1408	NAG	O5-C1	2.58	1.48	1.43
5	B	1403	NAG	O5-C5	2.51	1.48	1.43
5	A	1402	NAG	O5-C5	2.41	1.48	1.43
5	A	1409	NAG	C1-C2	2.34	1.55	1.52
5	B	1408	NAG	C1-C2	2.24	1.55	1.52
5	A	1402	NAG	O4-C4	2.23	1.48	1.43
5	C	1406	NAG	C1-C2	2.19	1.55	1.52
5	A	1406	NAG	C1-C2	2.09	1.55	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1401	NAG	C1-O5-C5	5.49	119.54	112.19
5	C	1402	NAG	C1-O5-C5	4.39	118.07	112.19
5	B	1403	NAG	C1-O5-C5	3.24	116.53	112.19
5	C	1403	NAG	C1-O5-C5	2.99	116.20	112.19
5	B	1407	NAG	O4-C4-C3	-2.72	103.96	110.38
5	A	1402	NAG	O5-C1-C2	-2.72	107.09	111.29
5	B	1404	NAG	C1-O5-C5	2.64	115.73	112.19
5	C	1407	NAG	O4-C4-C3	-2.34	104.86	110.38
5	B	1407	NAG	C1-O5-C5	2.19	115.12	112.19
5	A	1408	NAG	C1-O5-C5	2.18	115.10	112.19
5	A	1403	NAG	C1-O5-C5	2.17	115.09	112.19
5	B	1402	NAG	C1-O5-C5	2.17	115.09	112.19

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	1401	NAG	O5-C5-C6-O6
5	C	1401	NAG	C4-C5-C6-O6
5	B	1402	NAG	O5-C5-C6-O6
5	A	1403	NAG	O5-C5-C6-O6
5	B	1404	NAG	O5-C5-C6-O6
5	A	1405	NAG	C4-C5-C6-O6
5	A	1405	NAG	O5-C5-C6-O6
5	B	1401	NAG	C4-C5-C6-O6
5	A	1403	NAG	C4-C5-C6-O6
5	B	1402	NAG	C4-C5-C6-O6
5	C	1403	NAG	O5-C5-C6-O6
5	B	1404	NAG	C3-C2-N2-C7
5	C	1403	NAG	C3-C2-N2-C7

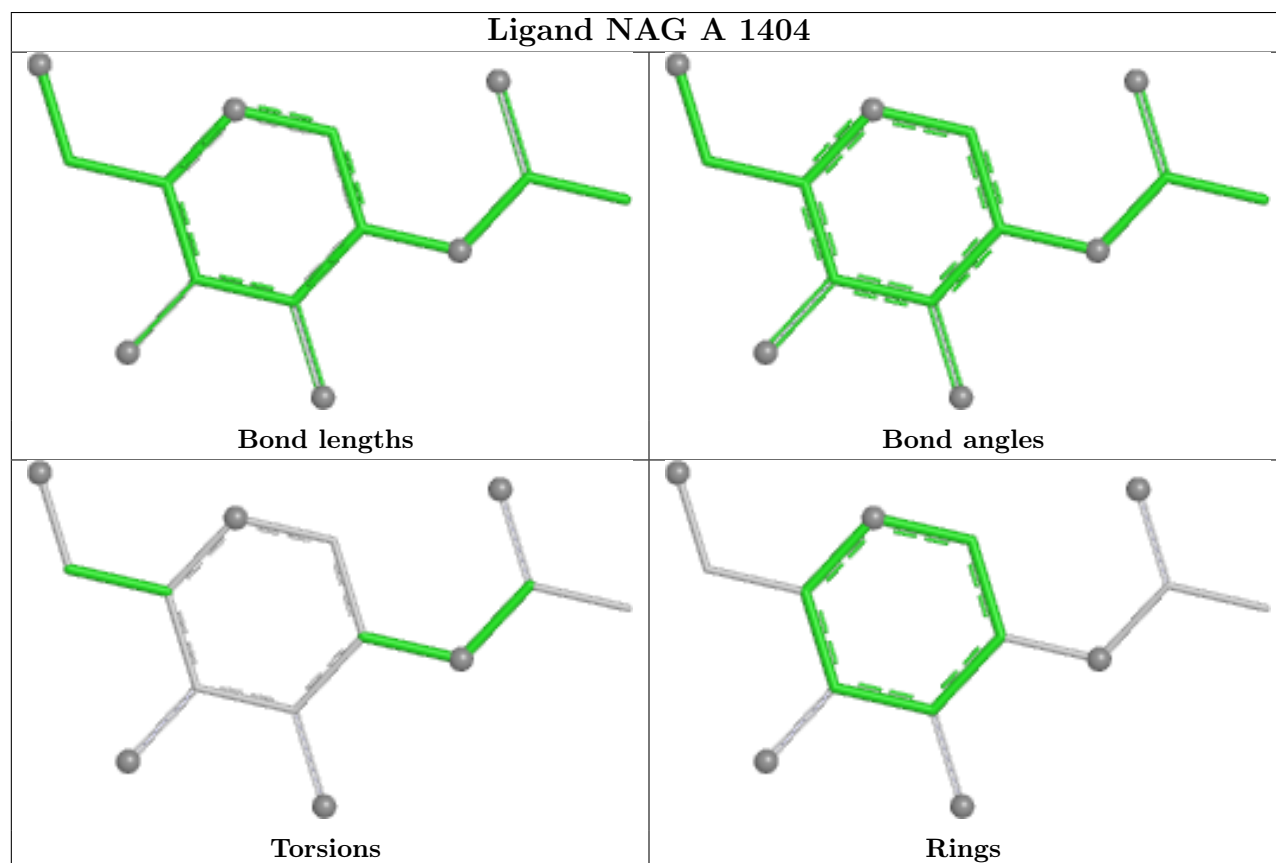
There are no ring outliers.

3 monomers are involved in 14 short contacts:

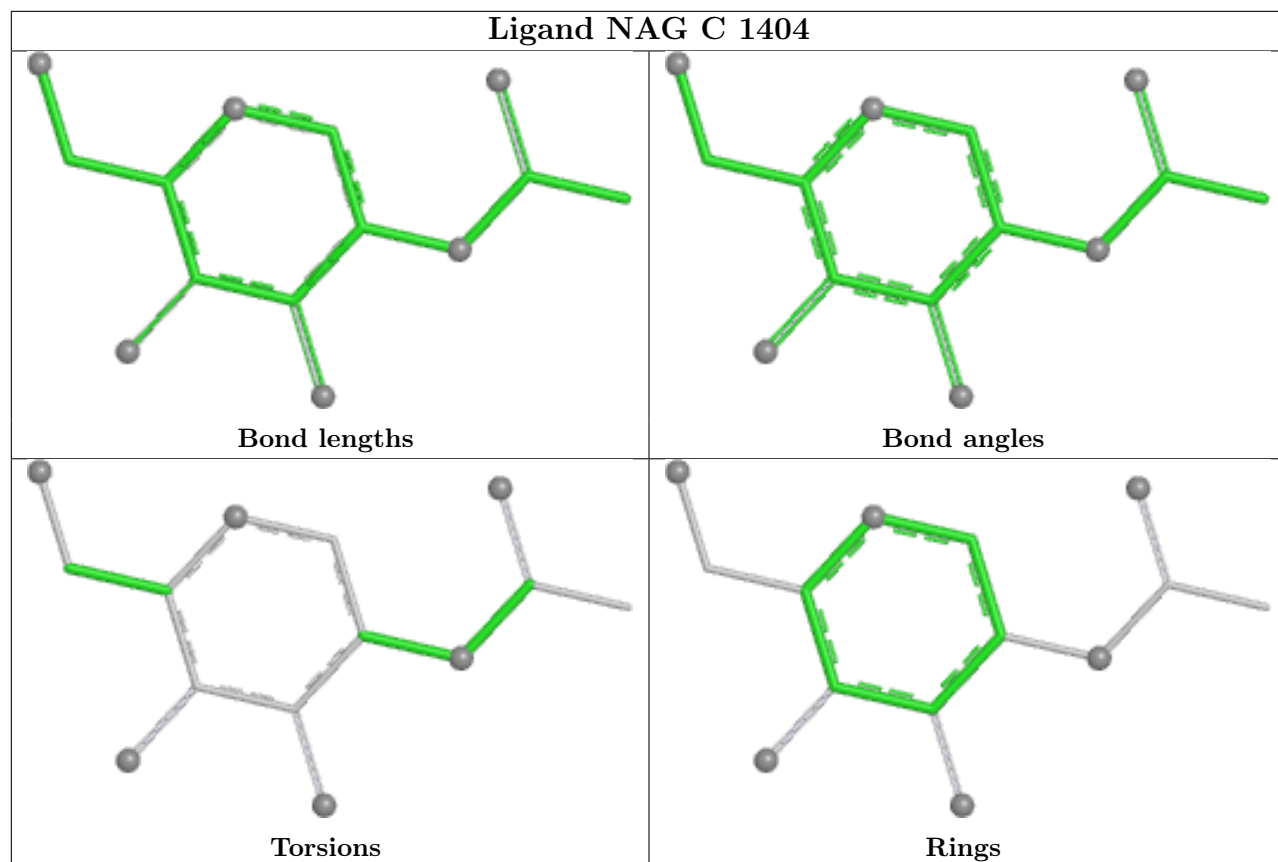
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1402	NAG	6	0
5	C	1406	NAG	2	0
5	B	1405	NAG	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

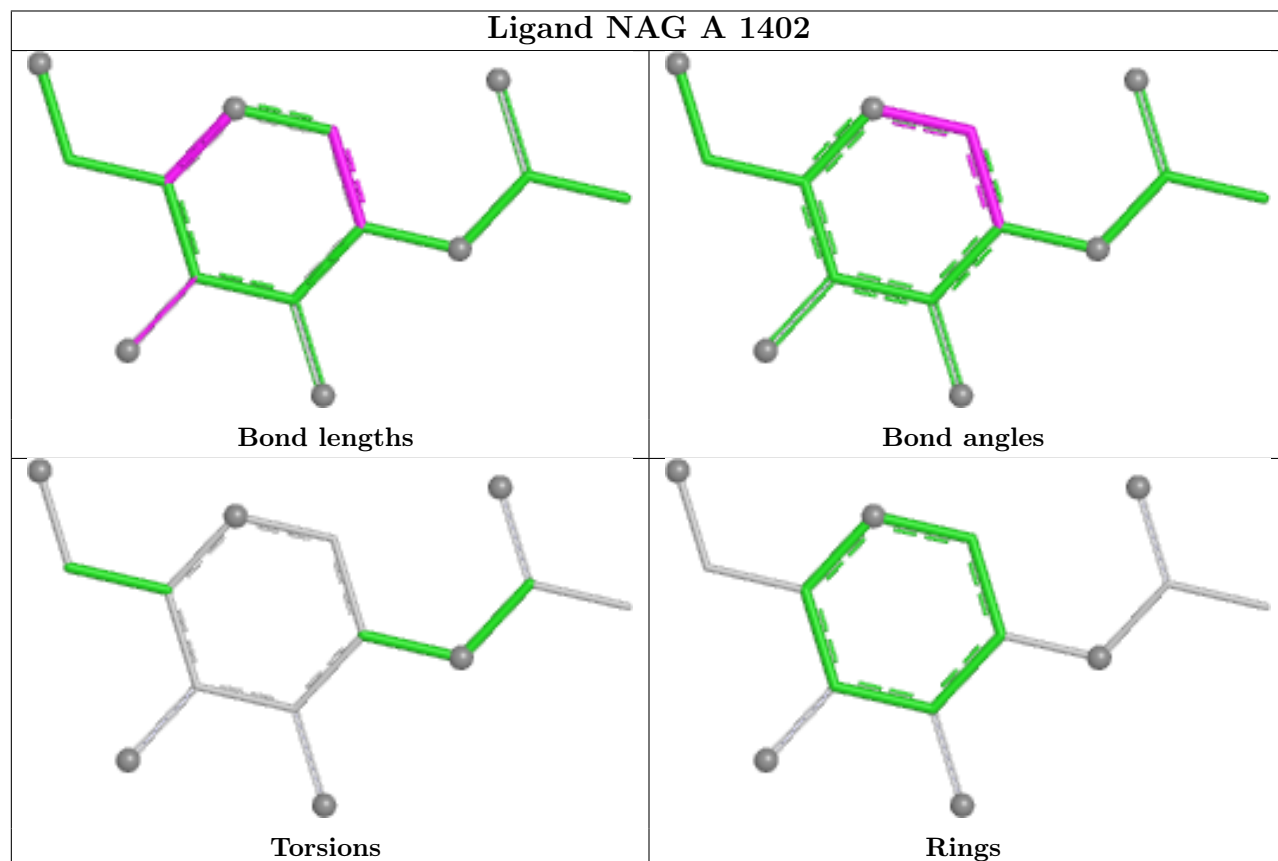
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



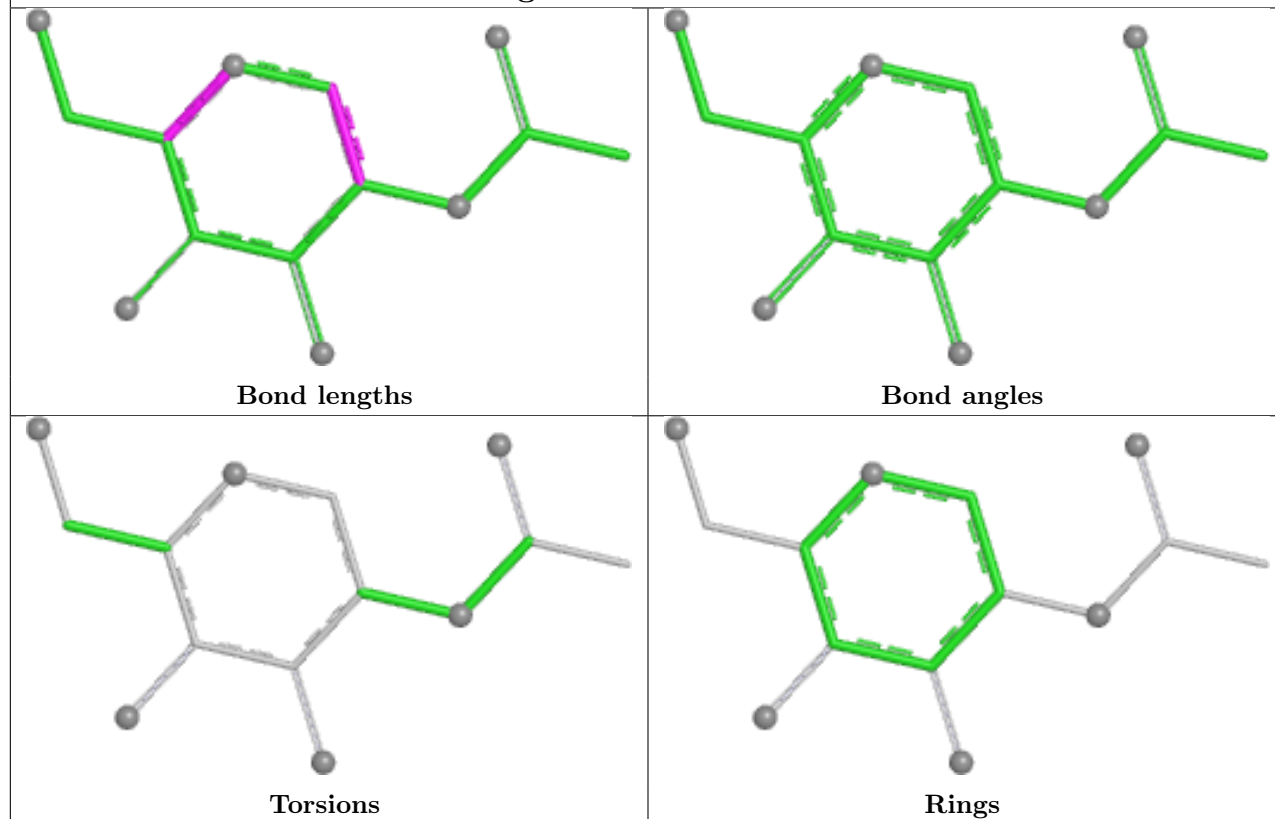
Ligand NAG C 1404



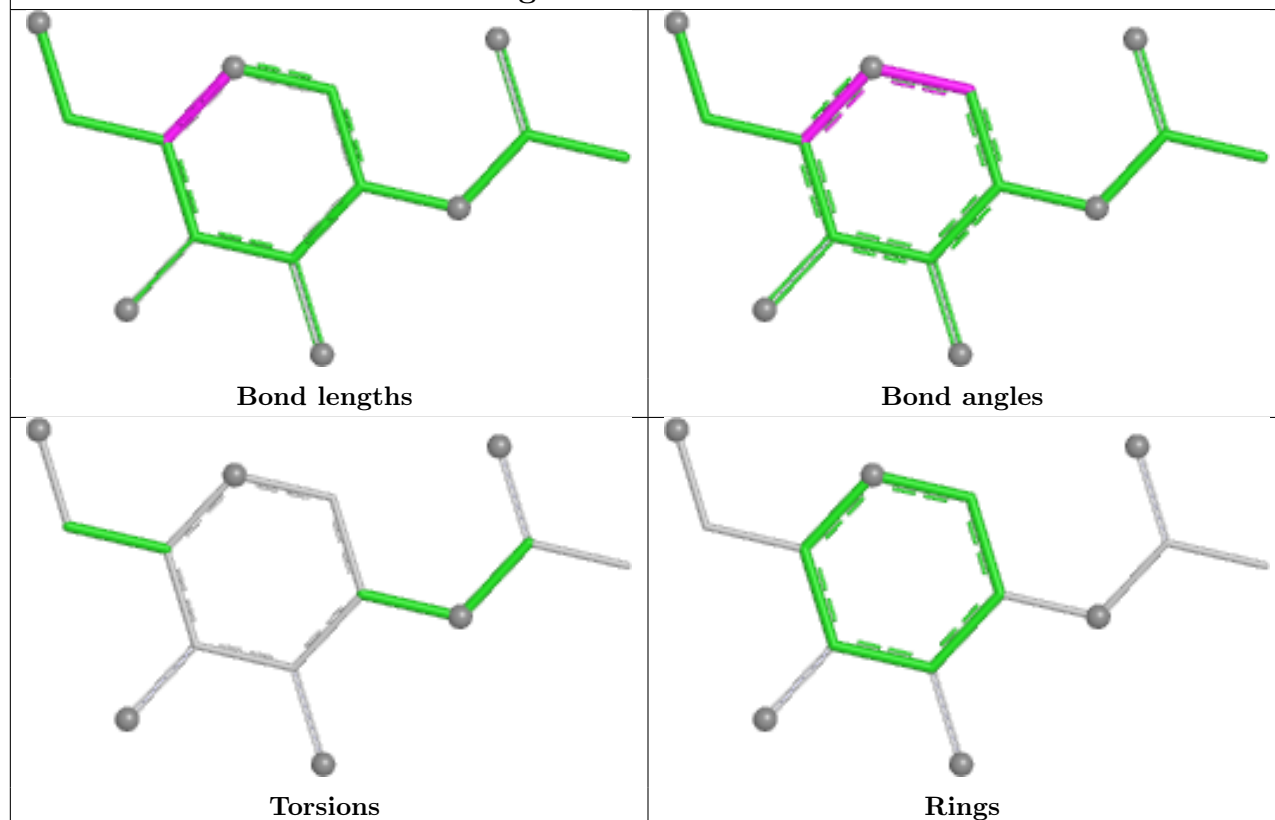
Ligand NAG A 1402



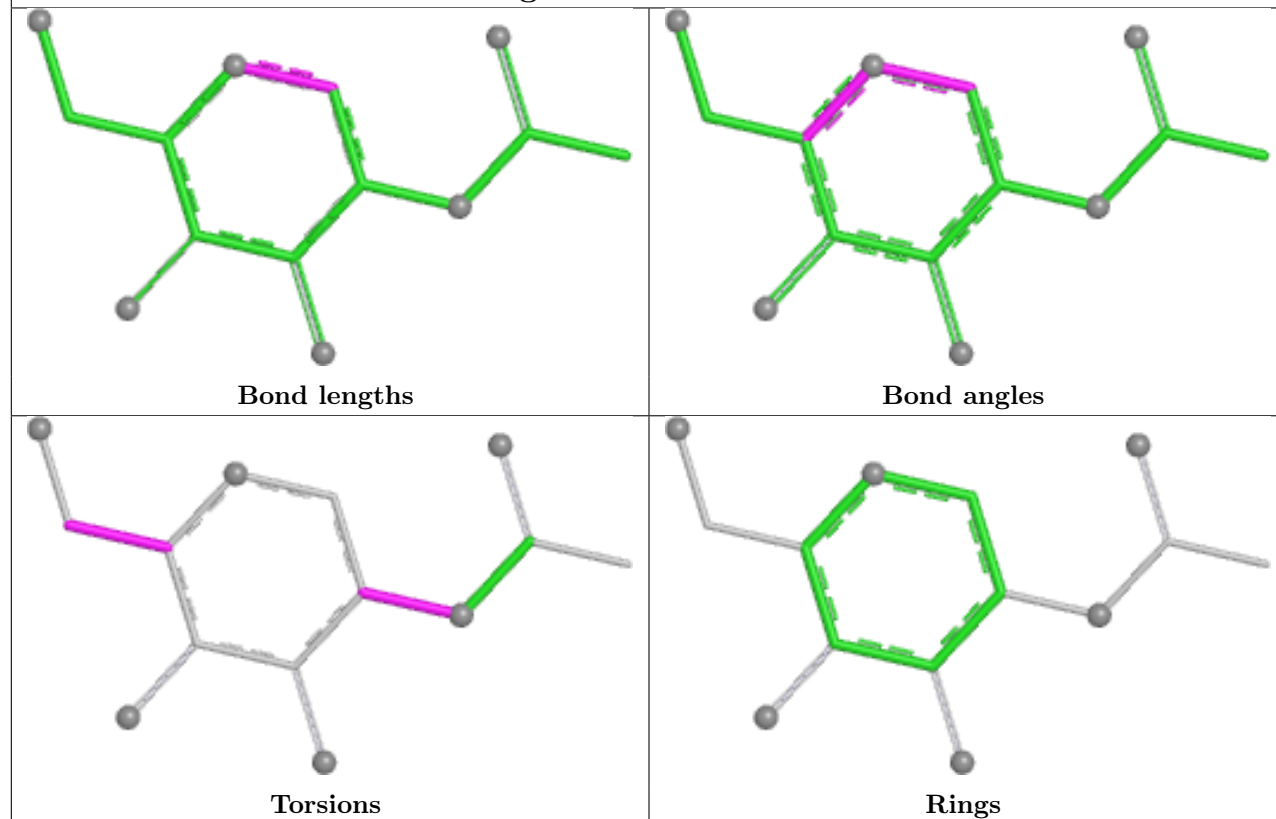
Ligand NAG A 1409



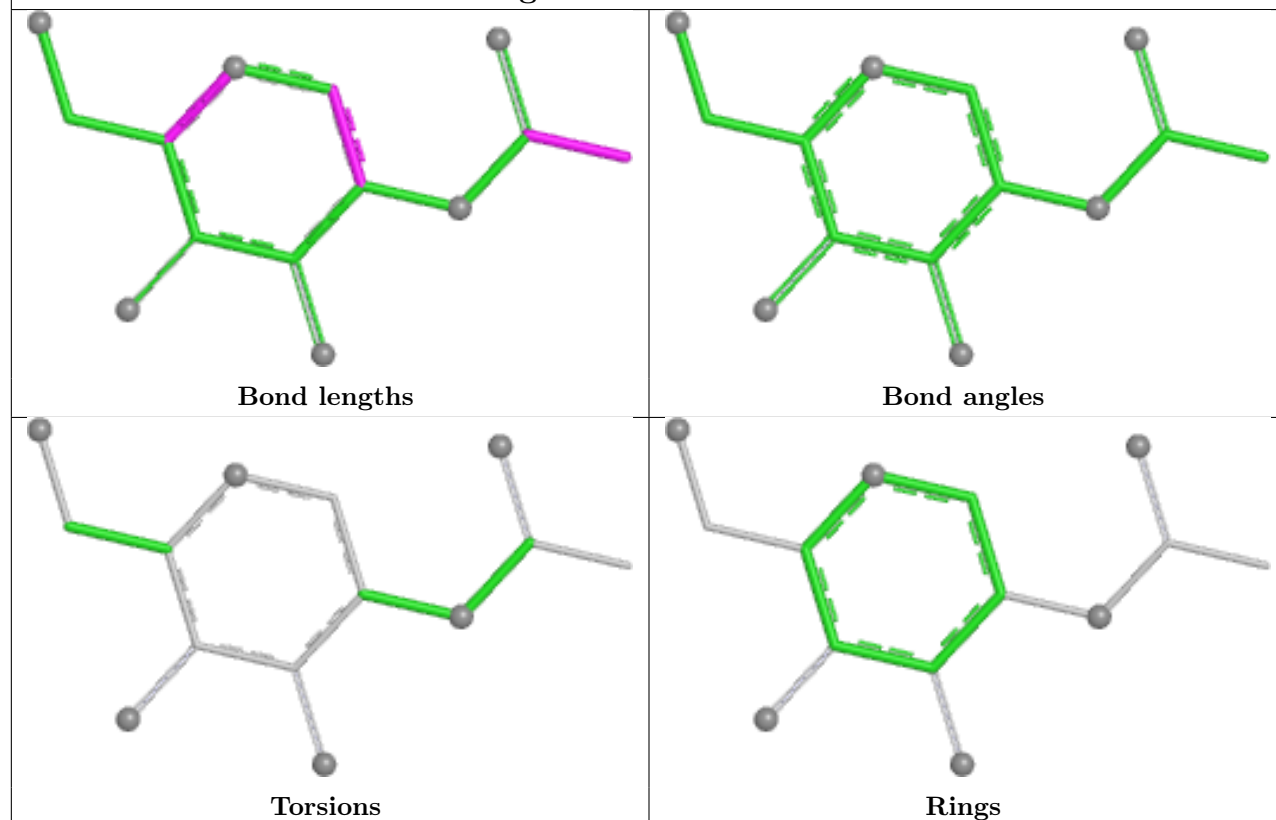
Ligand NAG B 1403



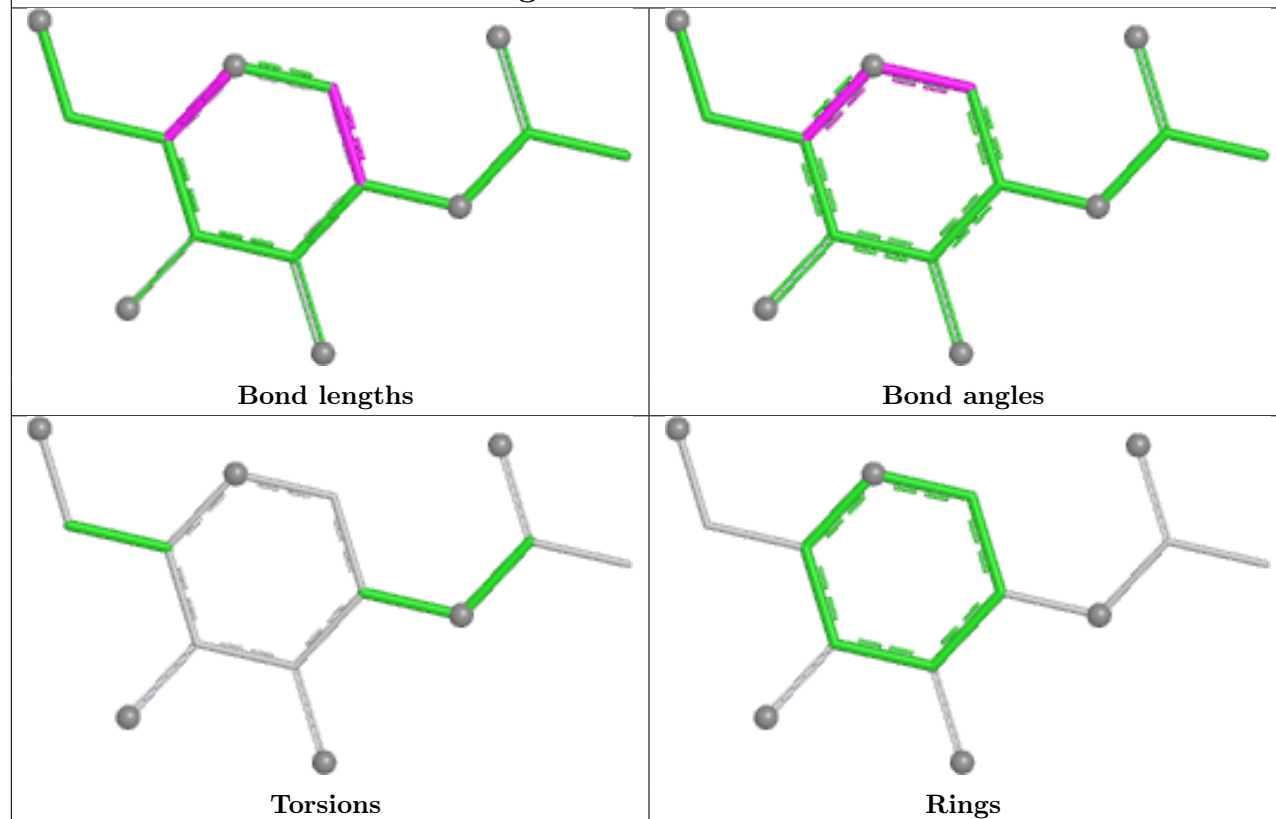
Ligand NAG B 1404



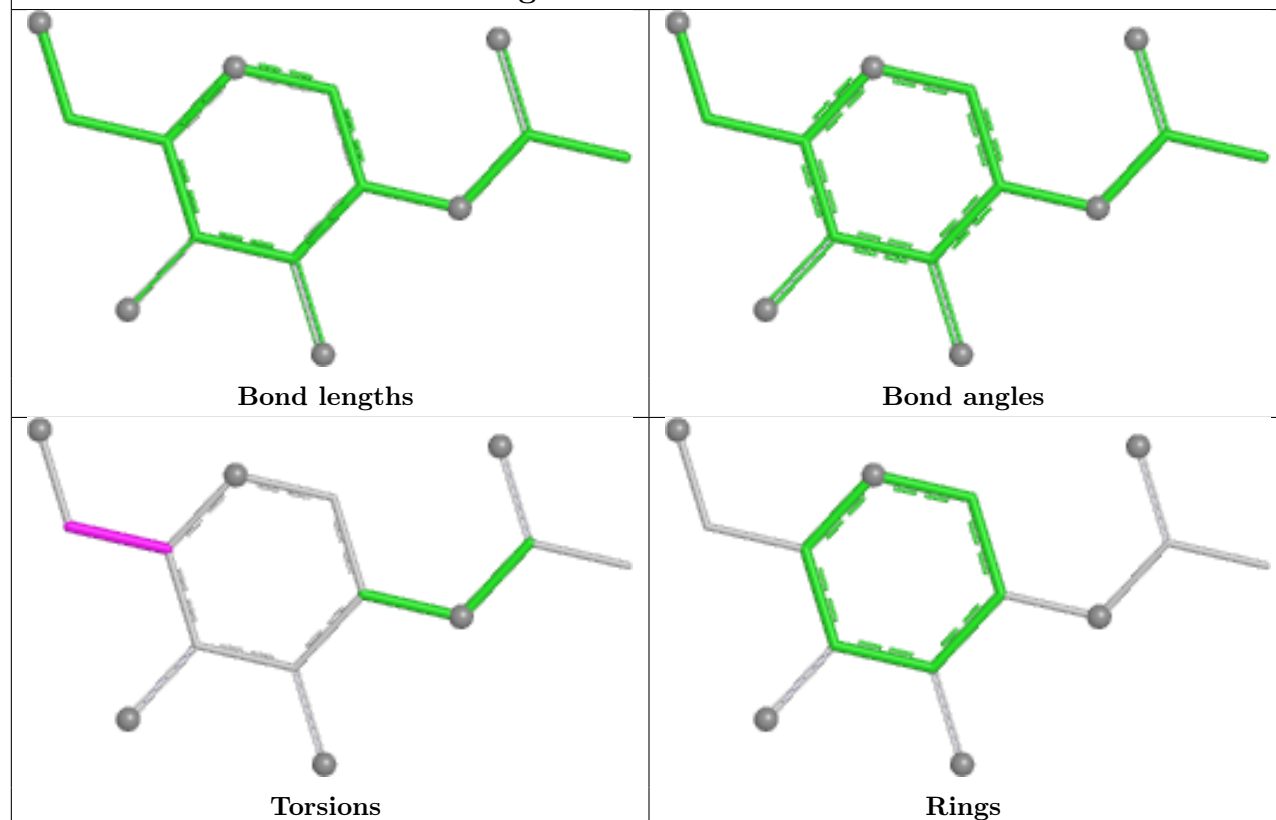
Ligand NAG C 1406

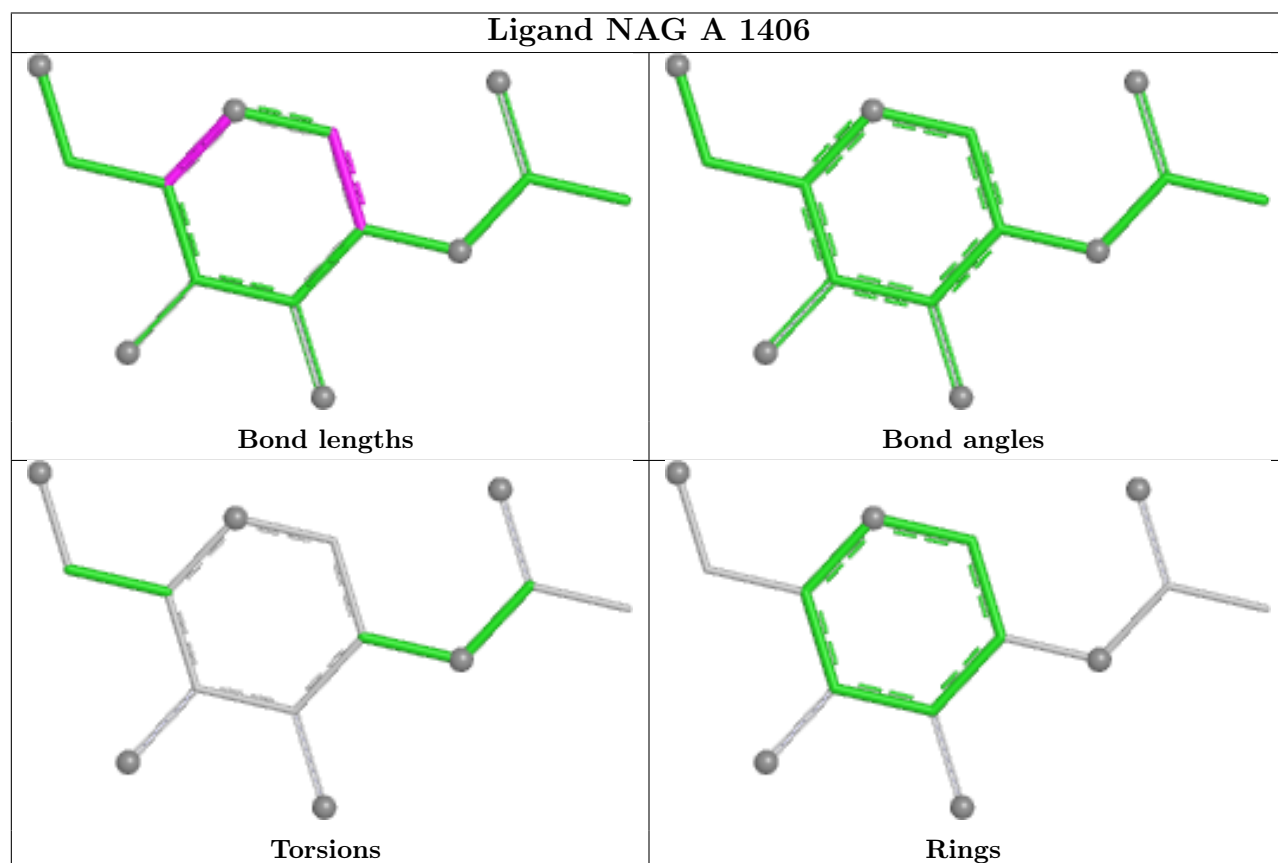
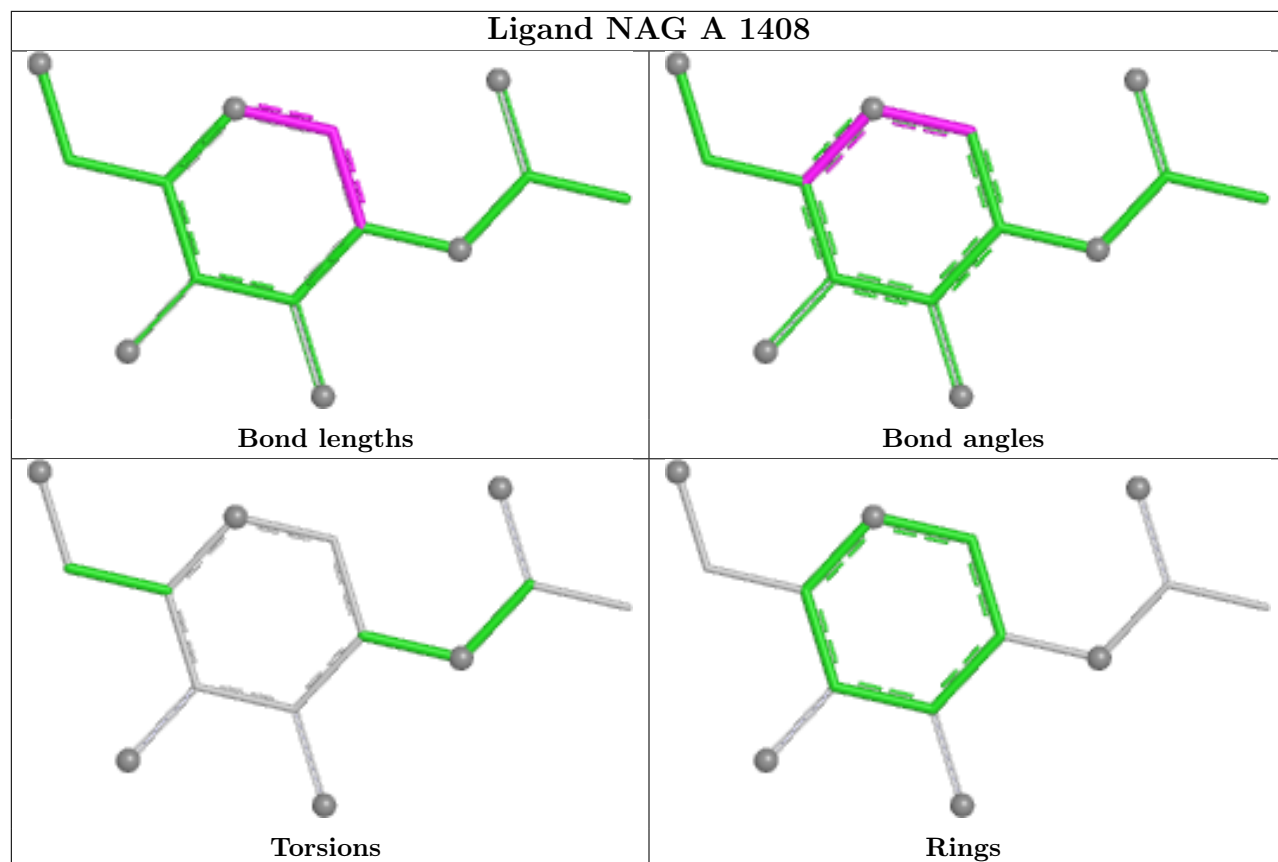


Ligand NAG A 1401

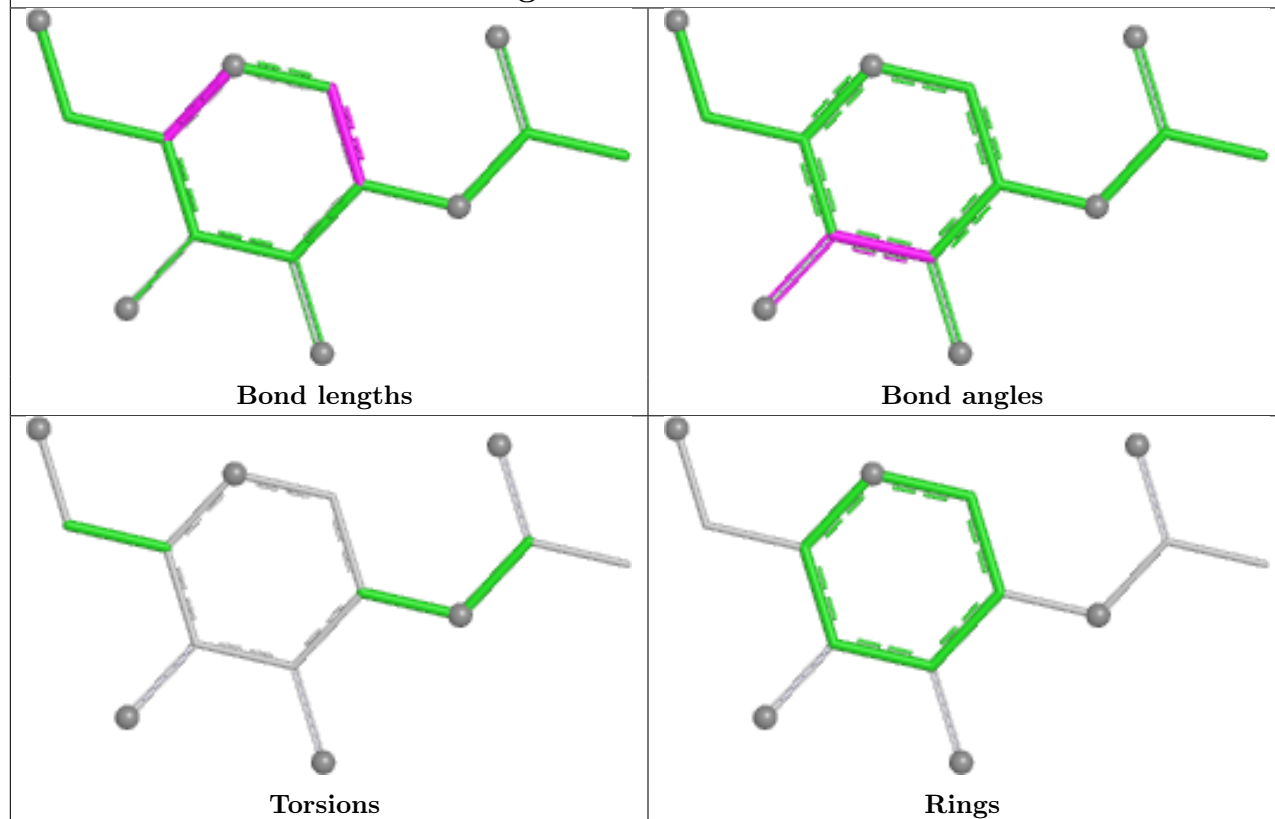


Ligand NAG A 1405

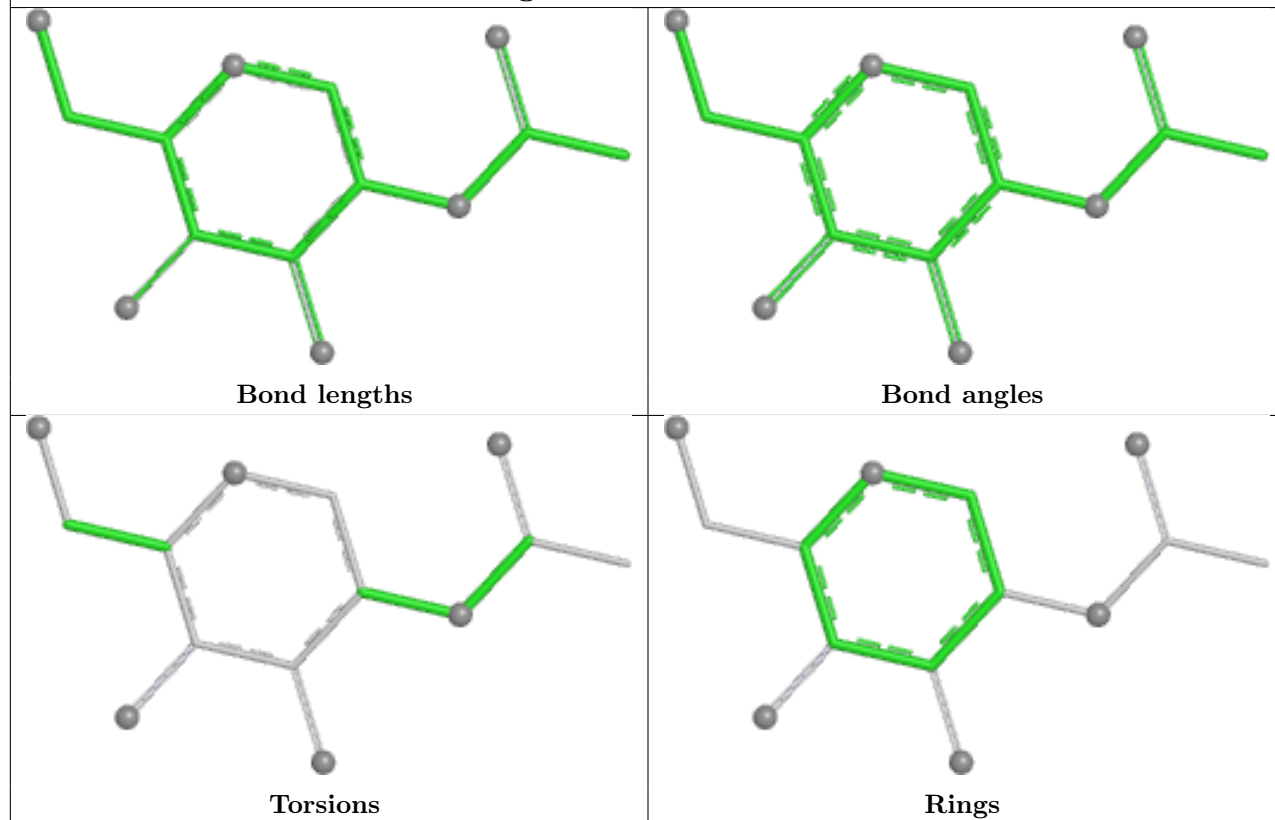


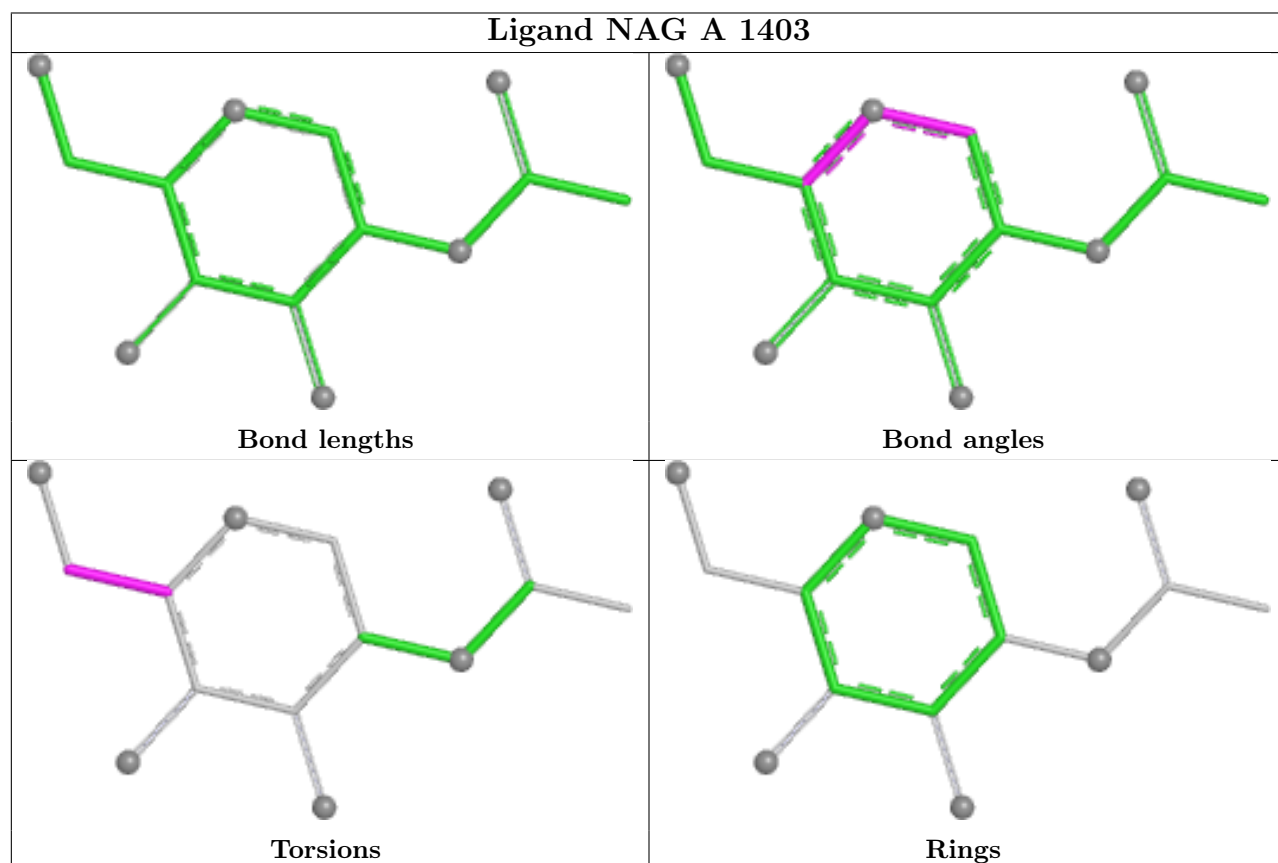
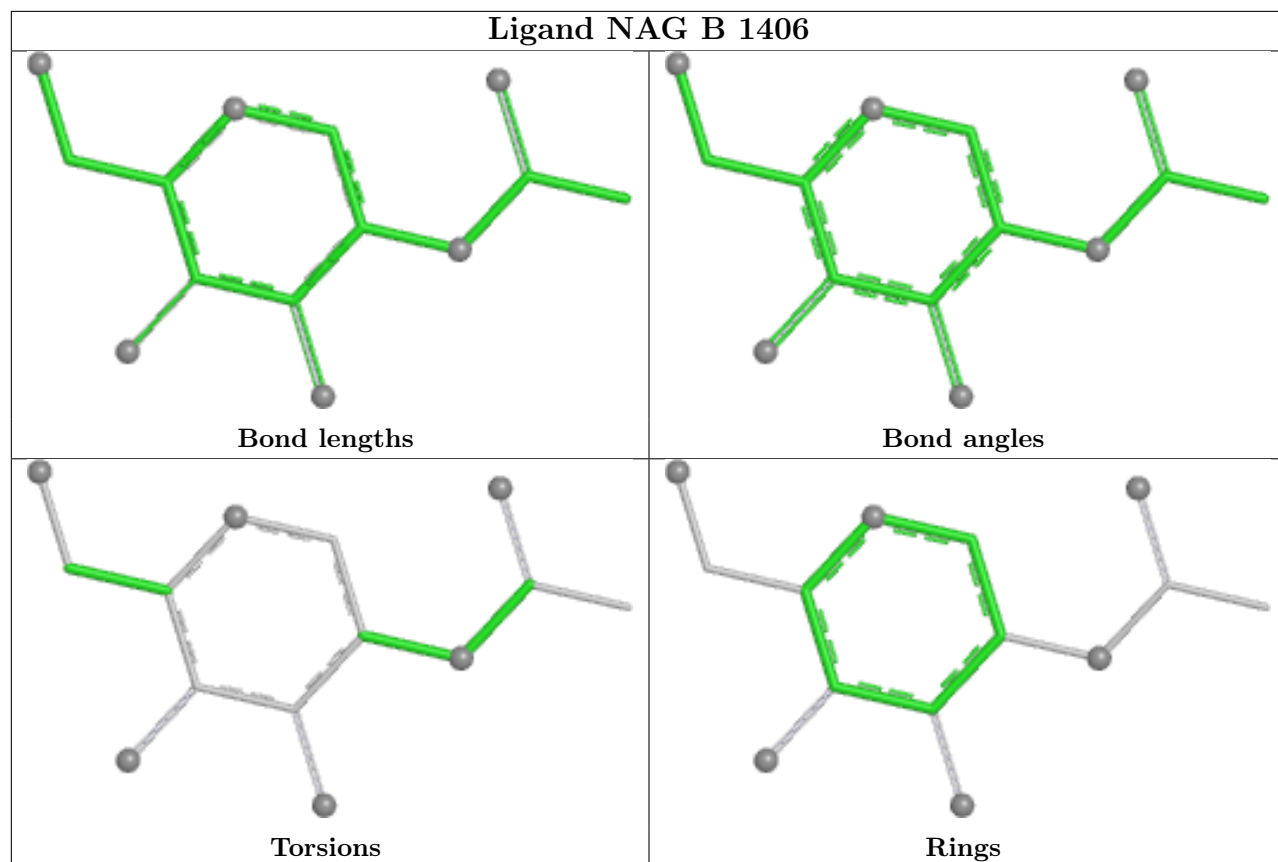


Ligand NAG C 1407

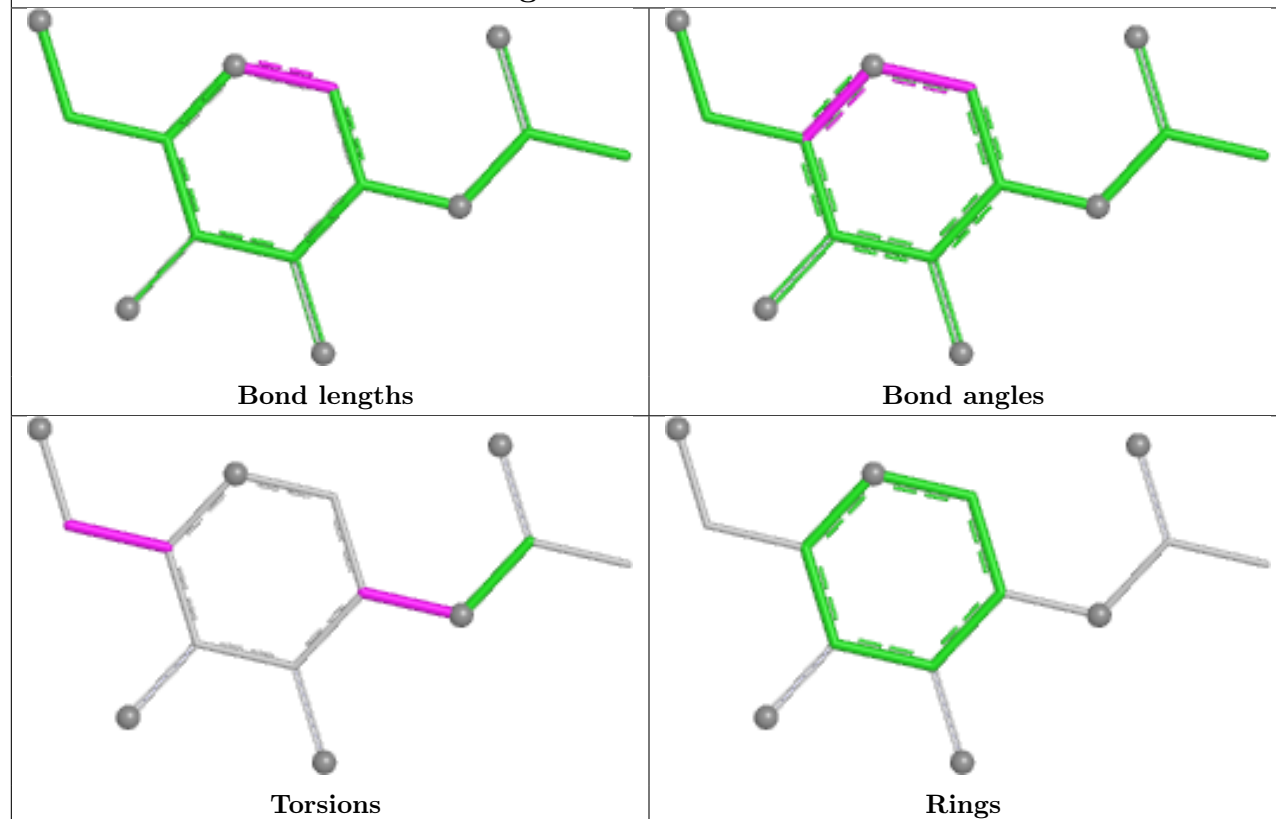


Ligand NAG A 1407

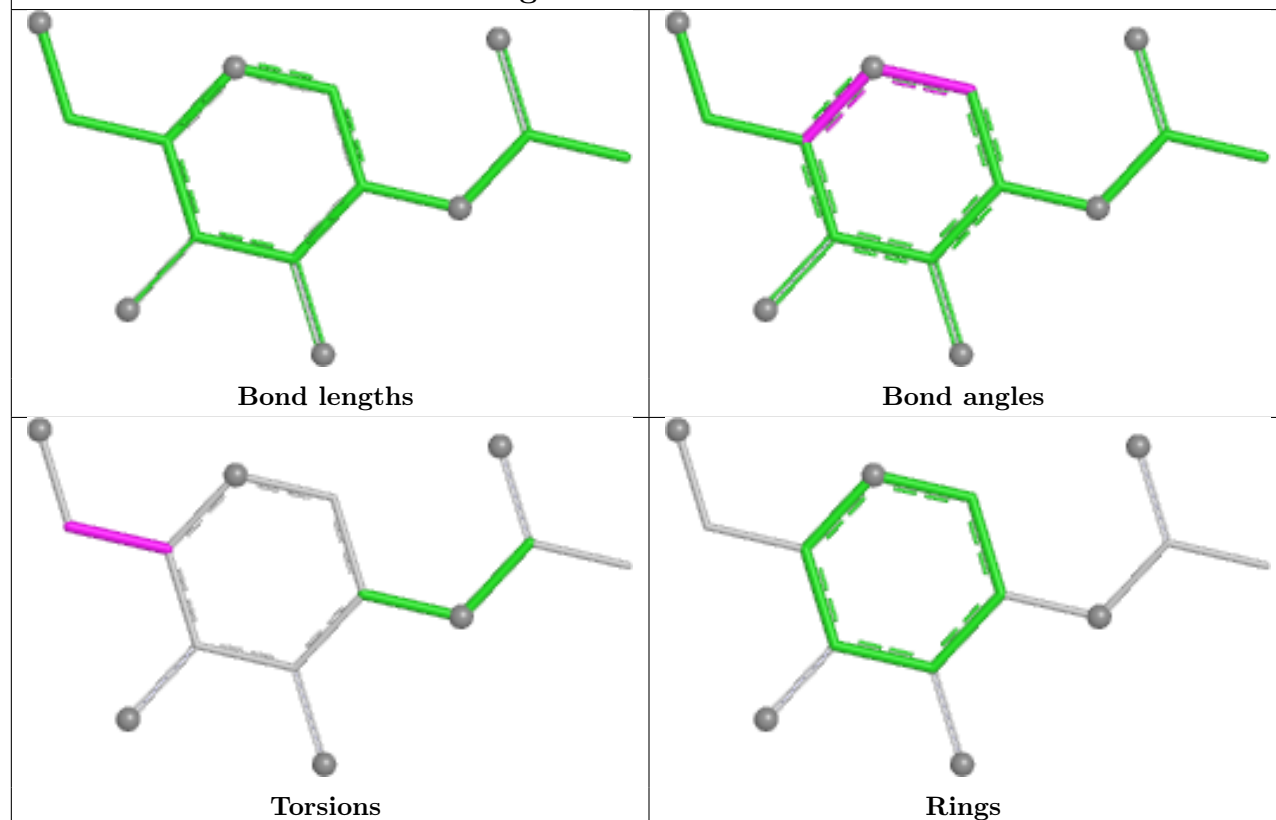


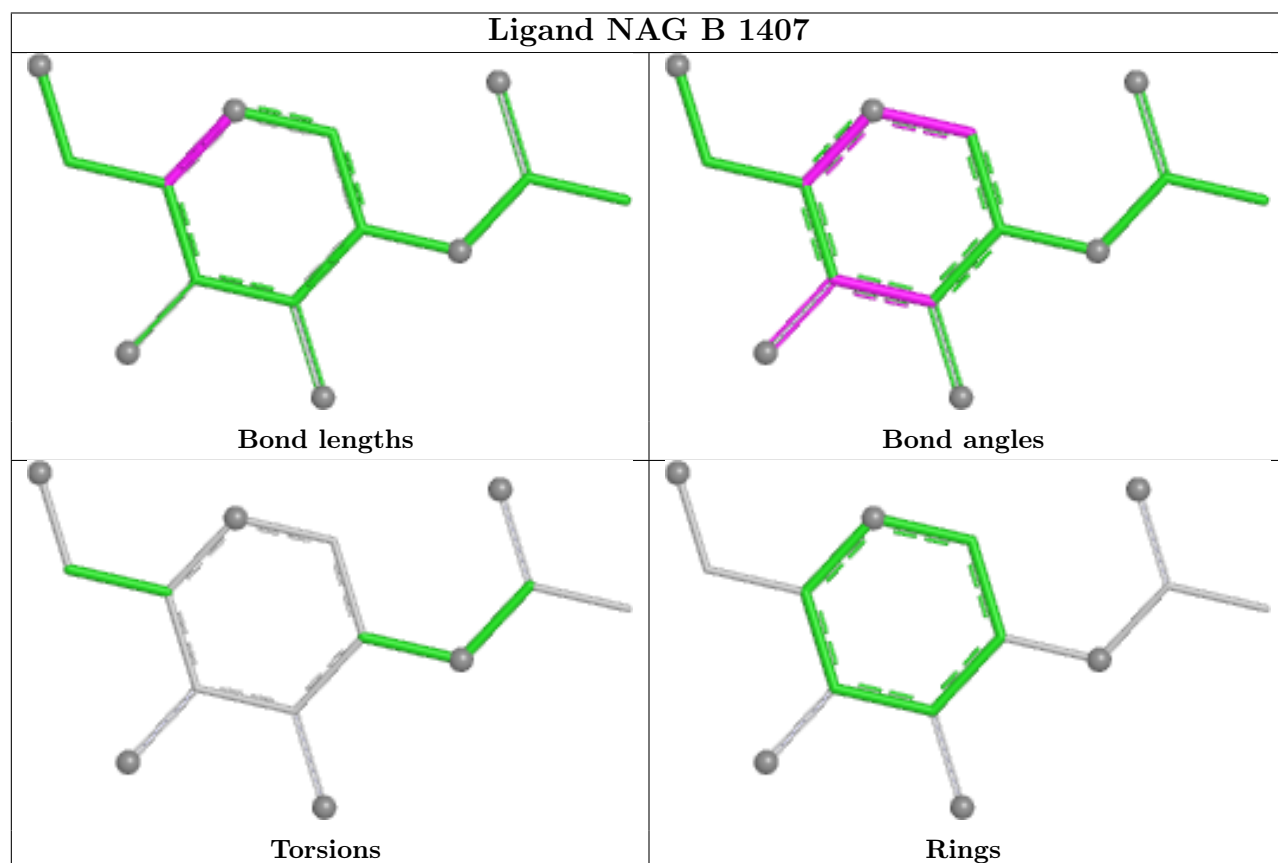
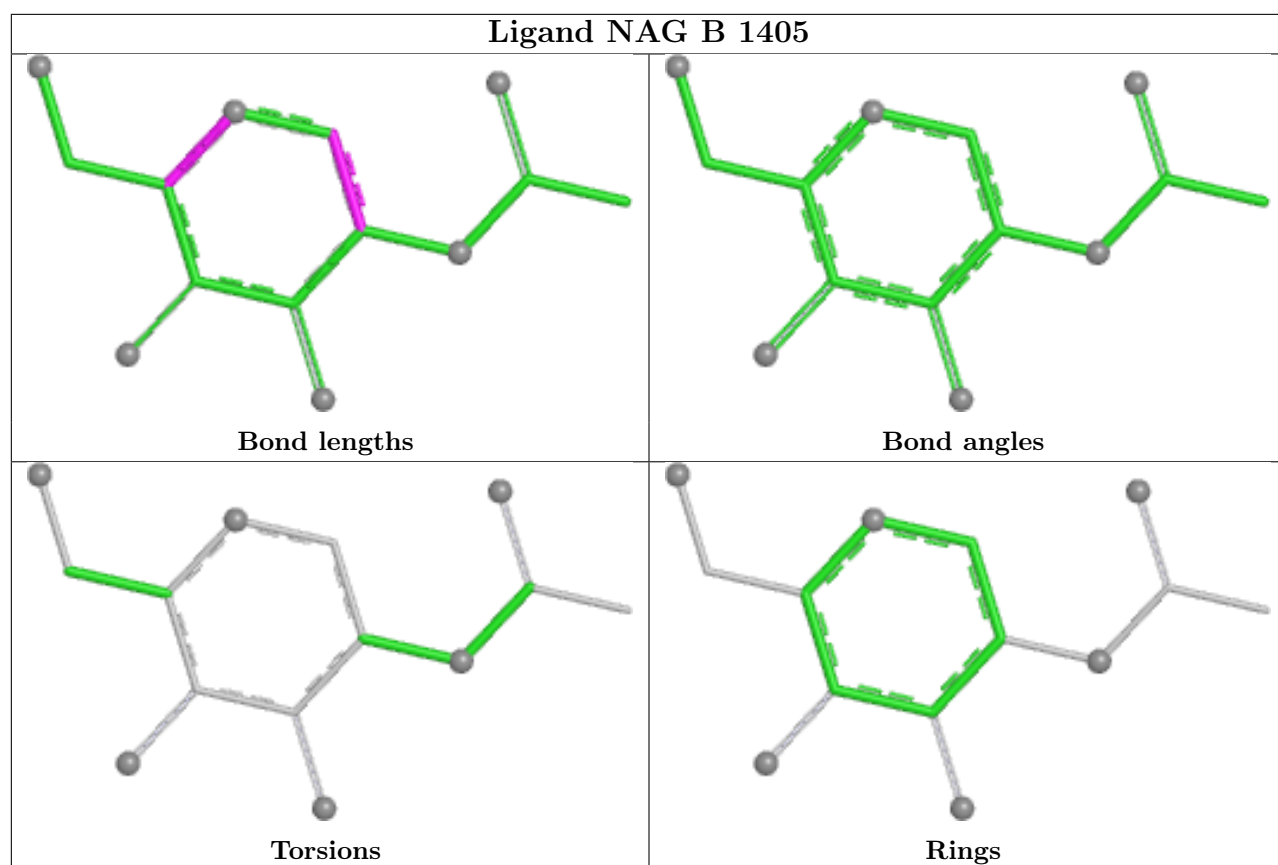


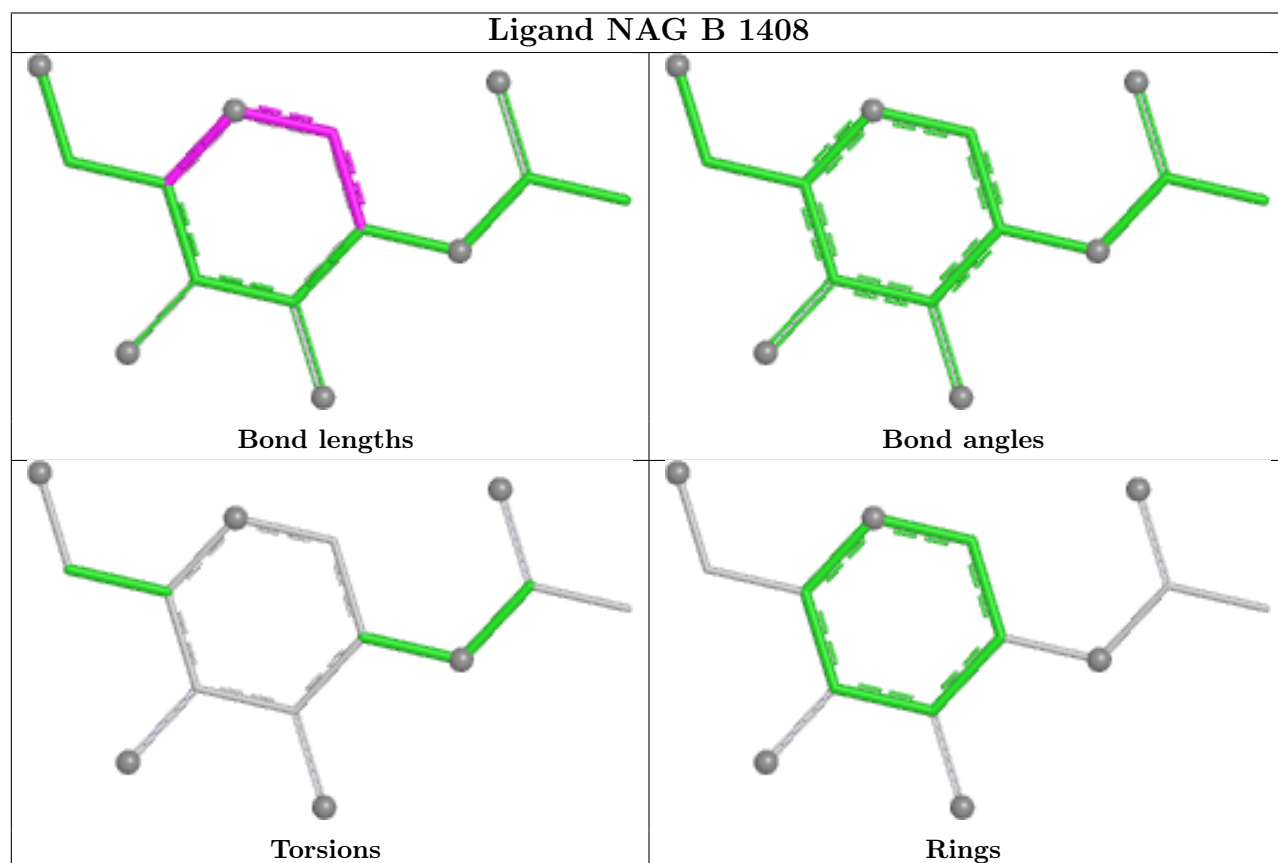
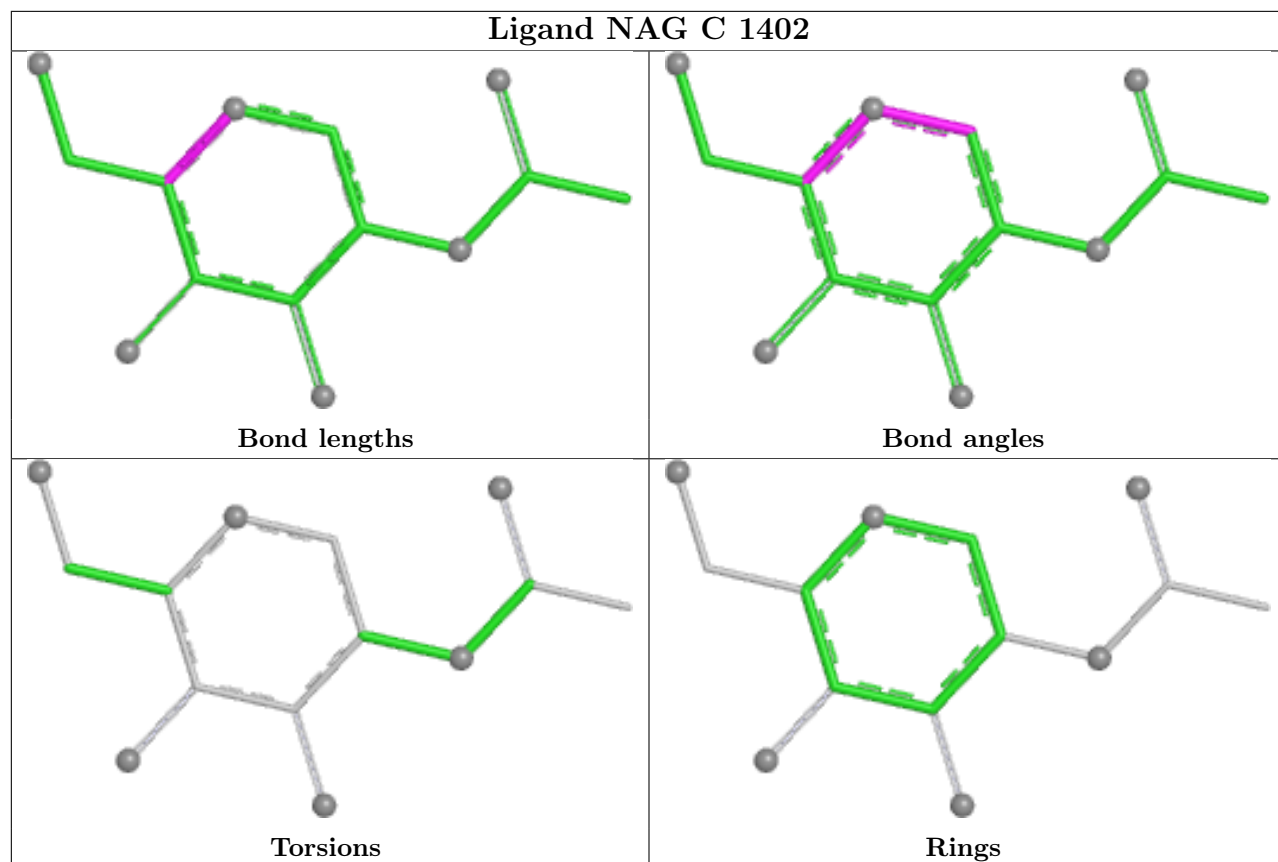
Ligand NAG C 1403

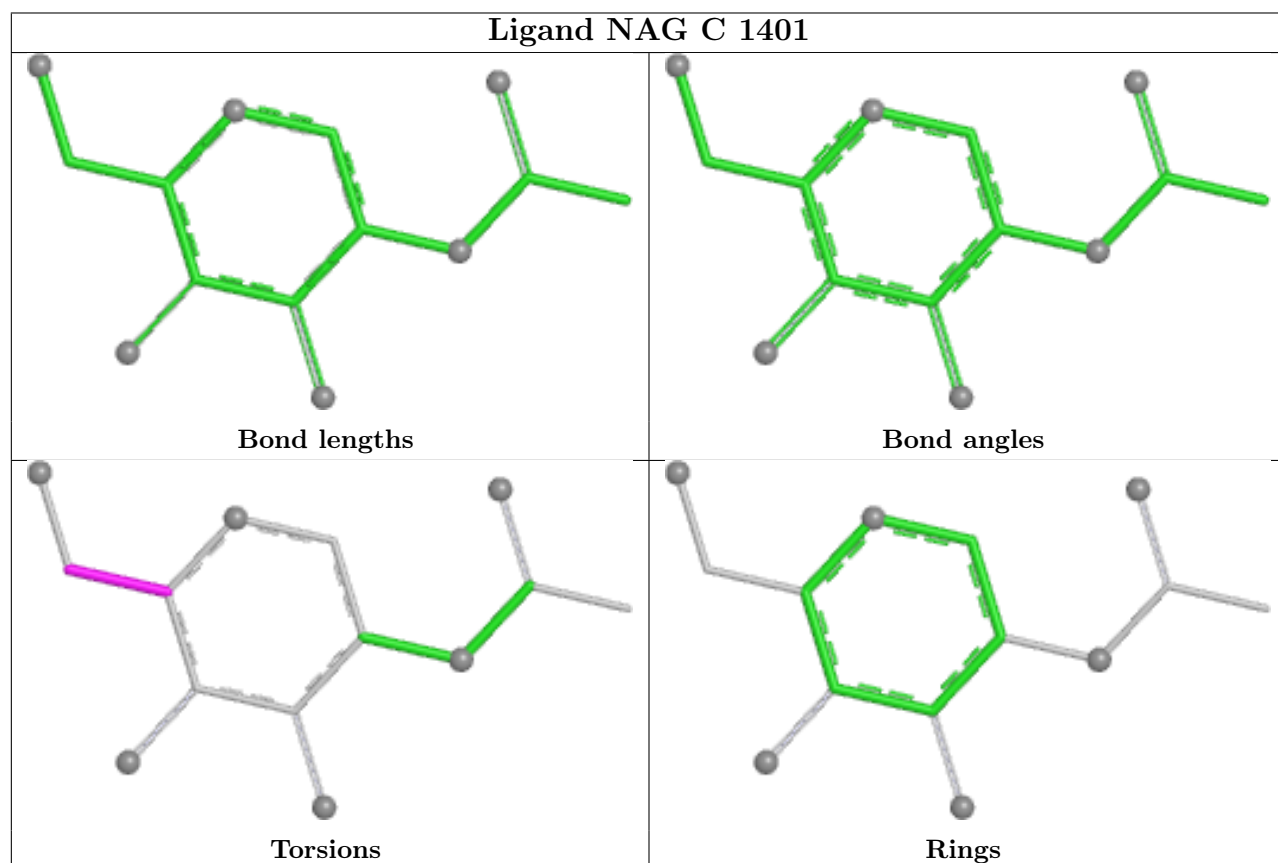
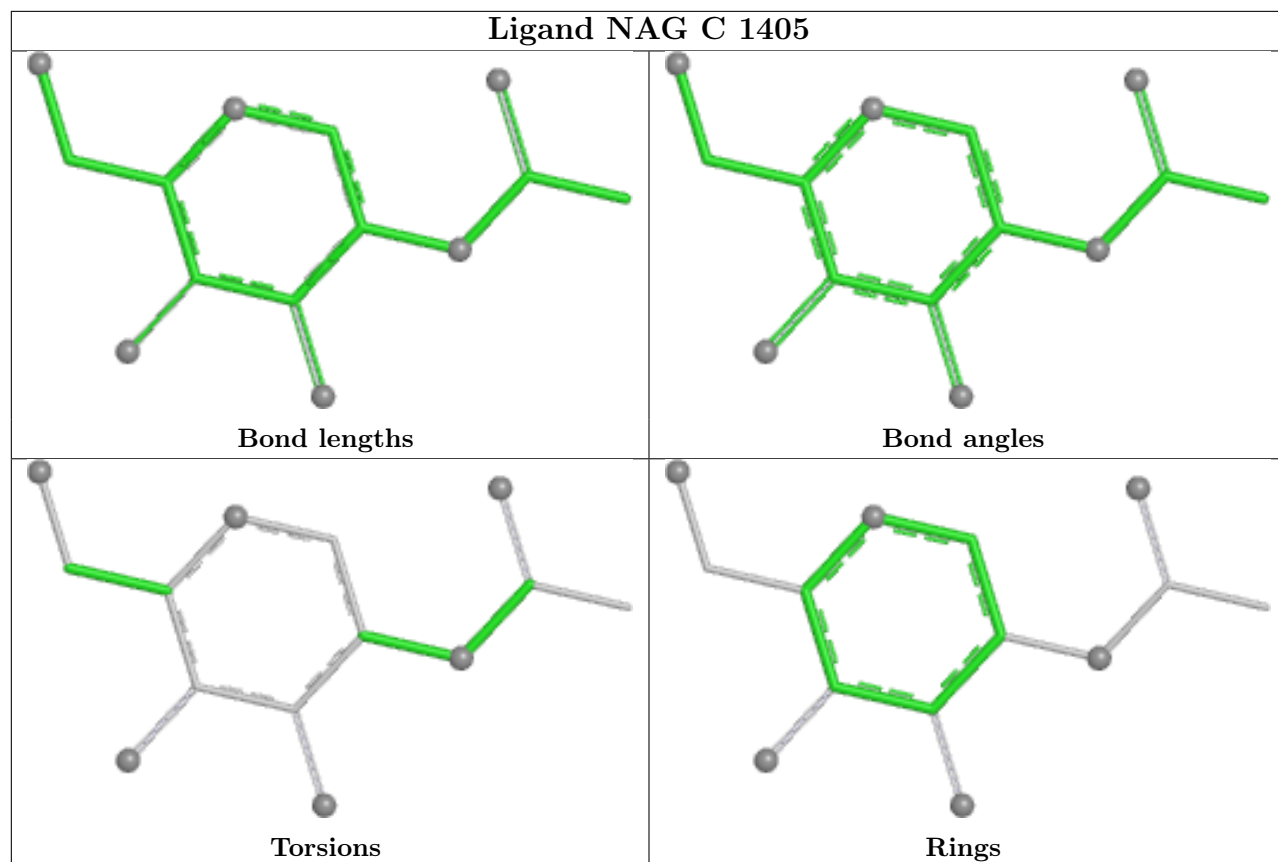


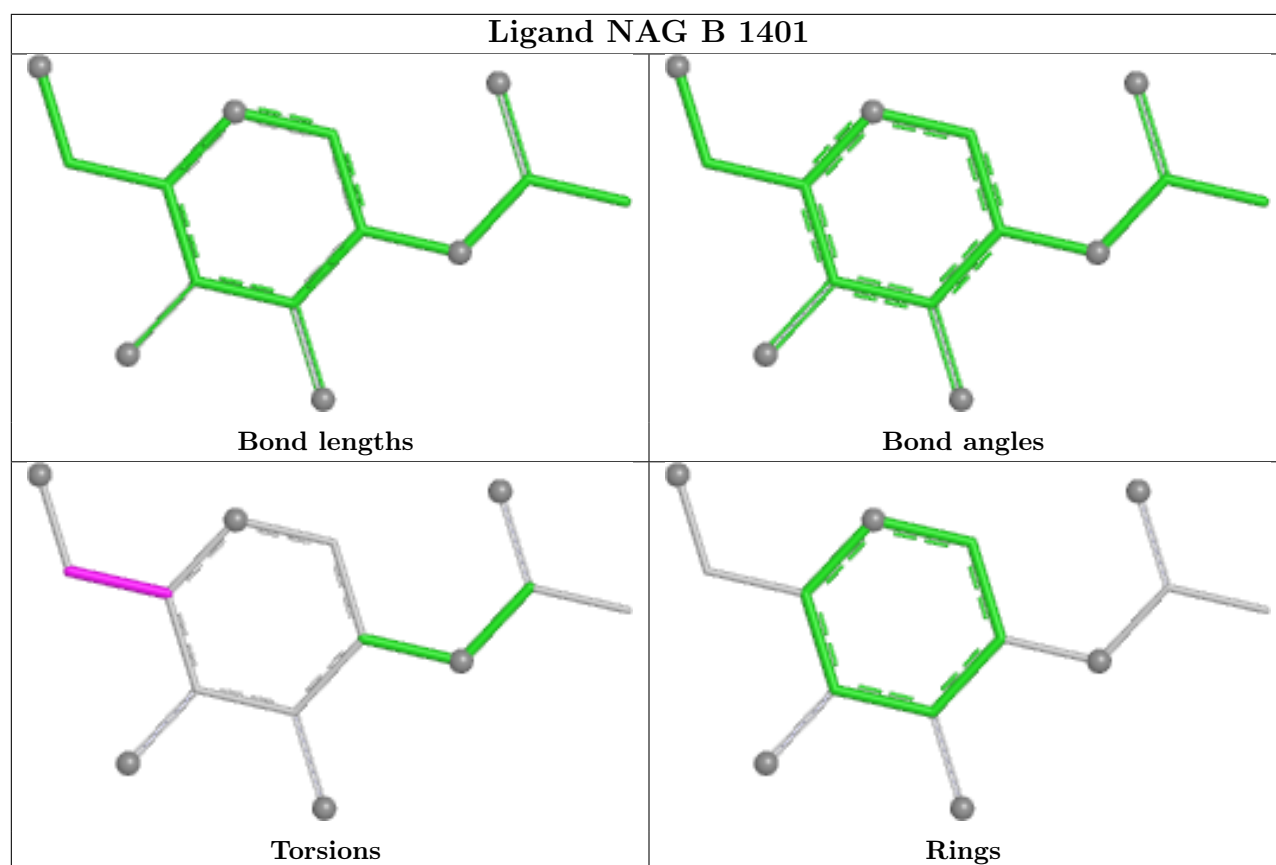
Ligand NAG B 1402











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.

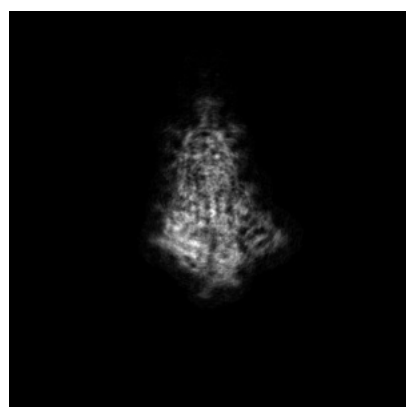
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24986. These allow visual inspection of the internal detail of the map and identification of artifacts.

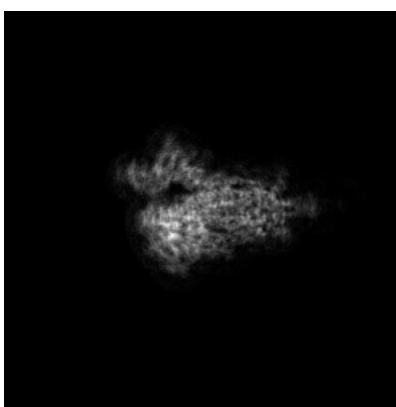
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

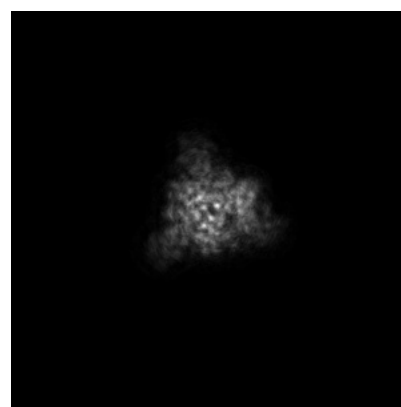
6.1.1 Primary map



X



Y



Z

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

6.2 Central slices [i](#)

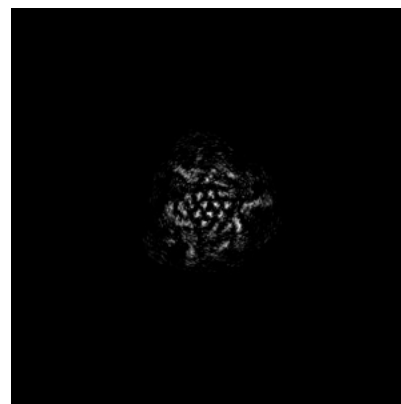
6.2.1 Primary map



X Index: 240



Y Index: 240



Z Index: 240

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

6.3 Largest variance slices [i](#)

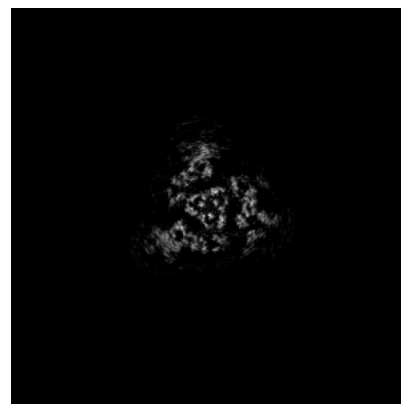
6.3.1 Primary map



X Index: 231



Y Index: 245

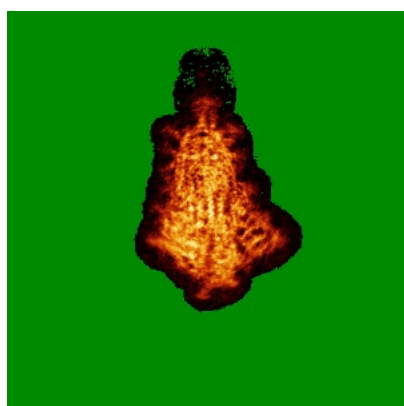


Z Index: 221

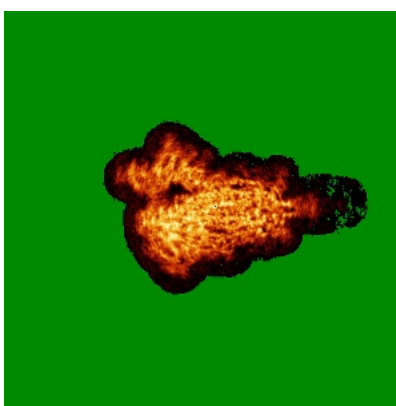
The images above show the largest variance slices of the map in three orthogonal directions.

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

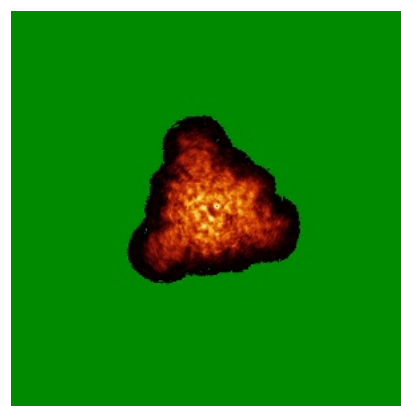
6.4.1 Primary map



X



Y

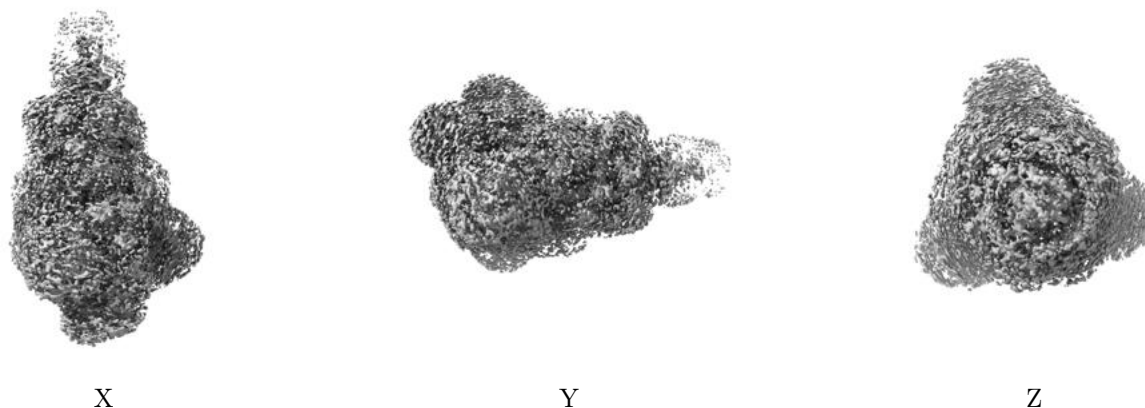


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

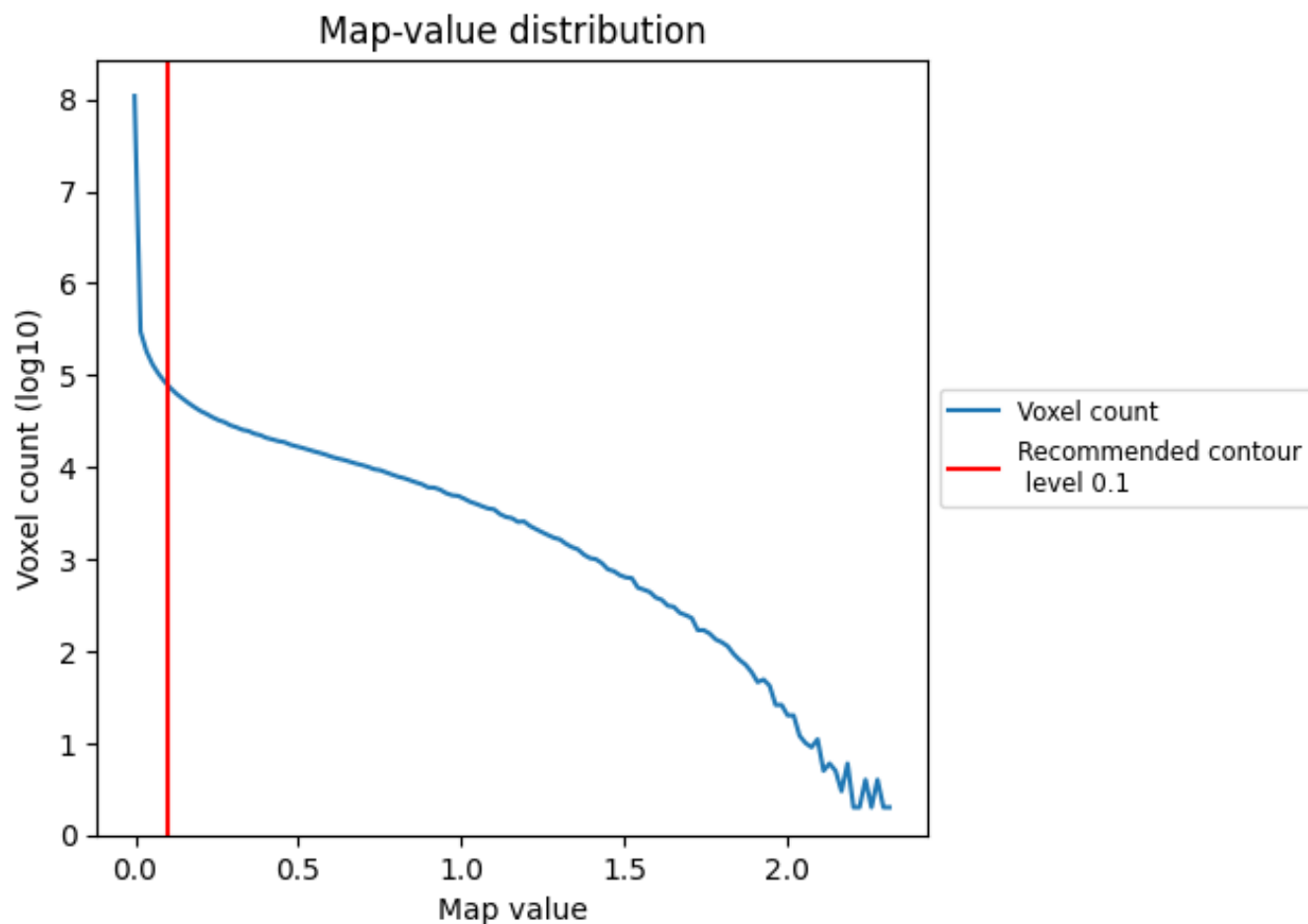
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

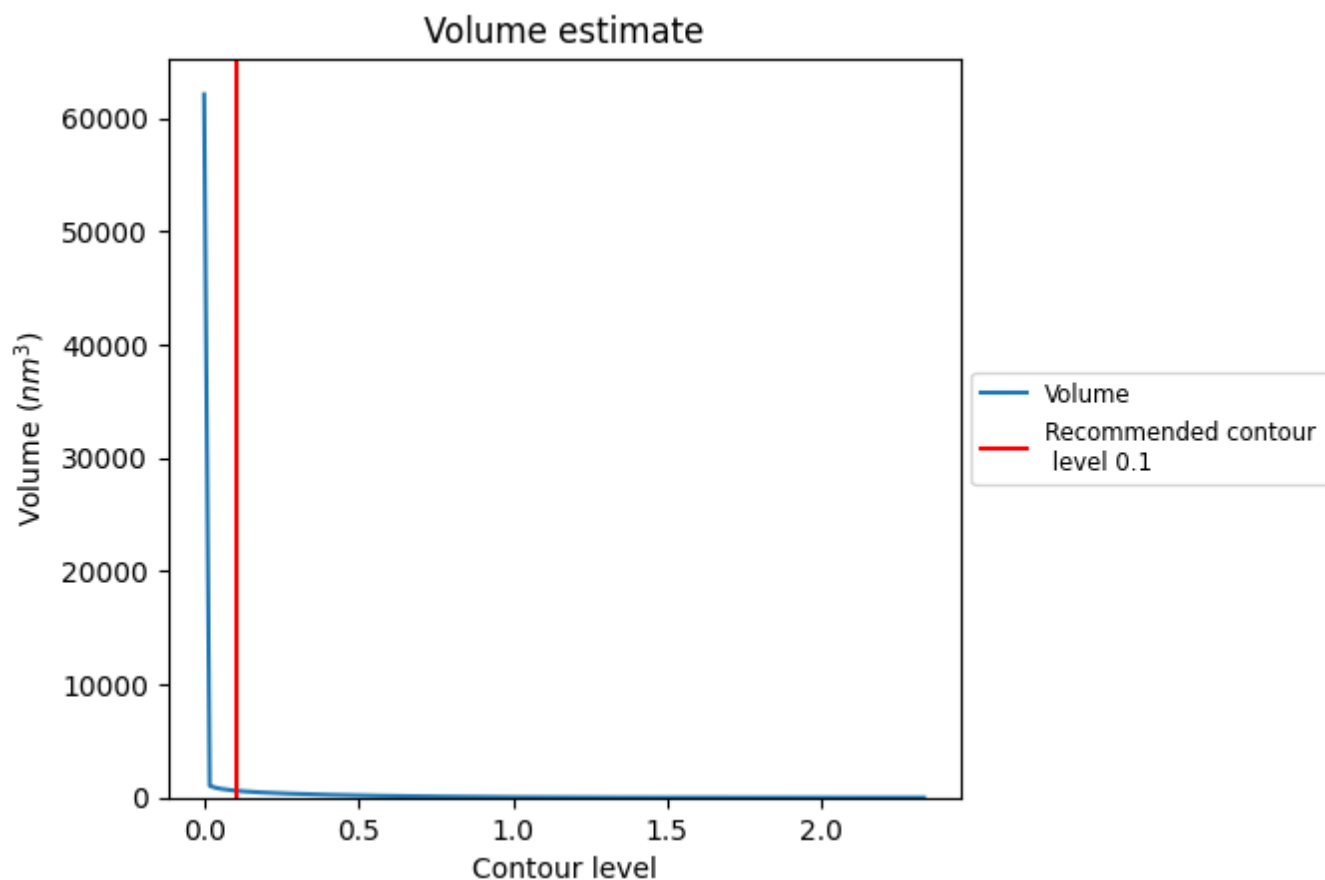
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

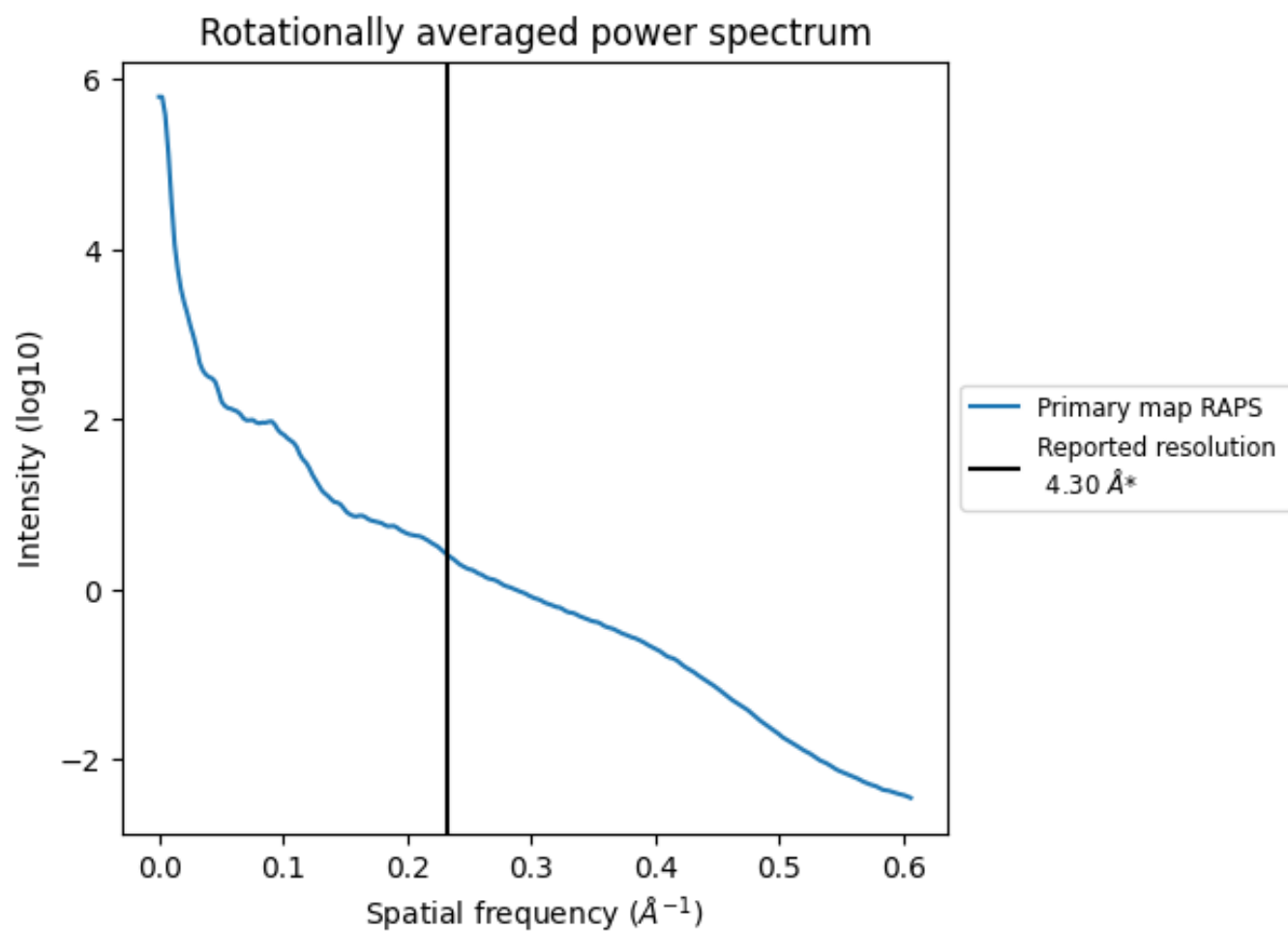
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 612 nm^3 ; this corresponds to an approximate mass of 553 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.233 Å⁻¹

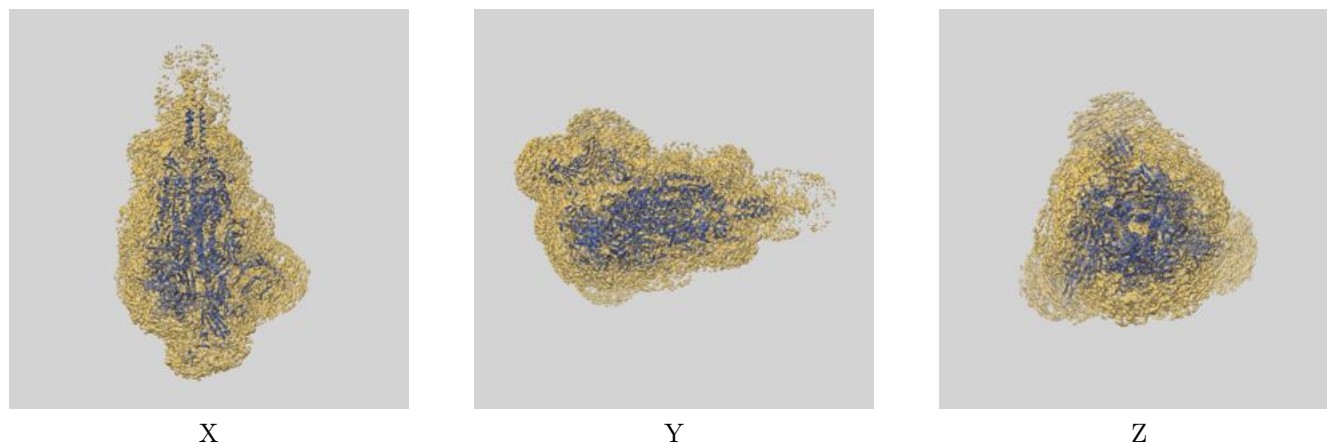
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

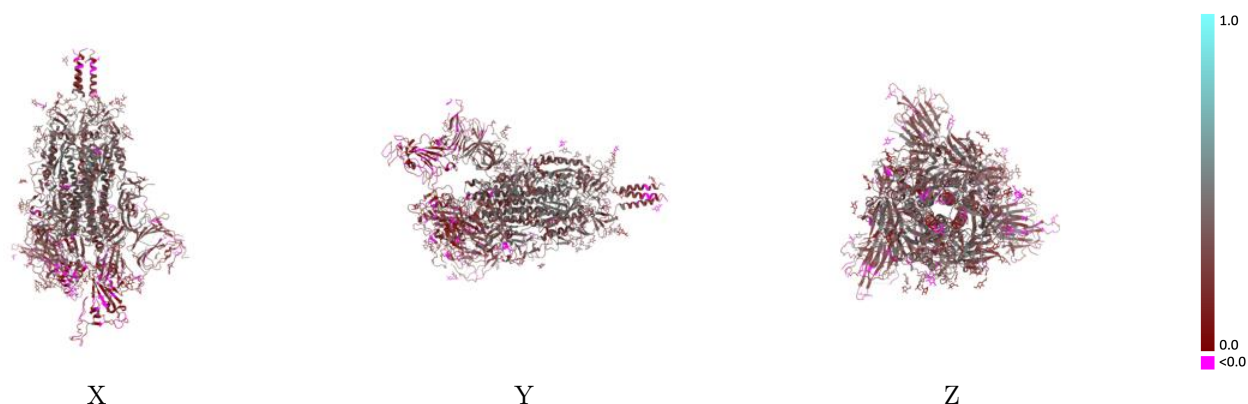
This section contains information regarding the fit between EMDB map EMD-24986 and PDB model 7SBR. Per-residue inclusion information can be found in section [3](#) on page [13](#).

9.1 Map-model overlay [i](#)



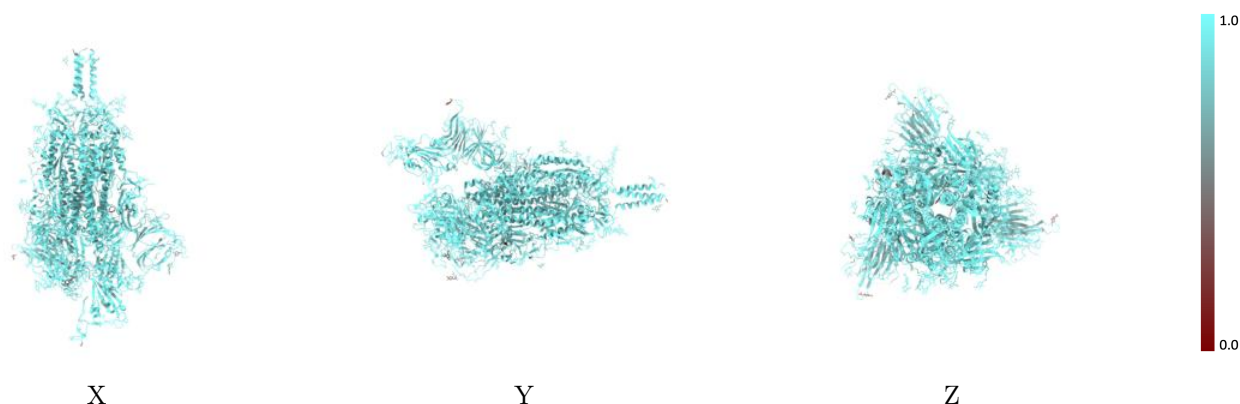
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



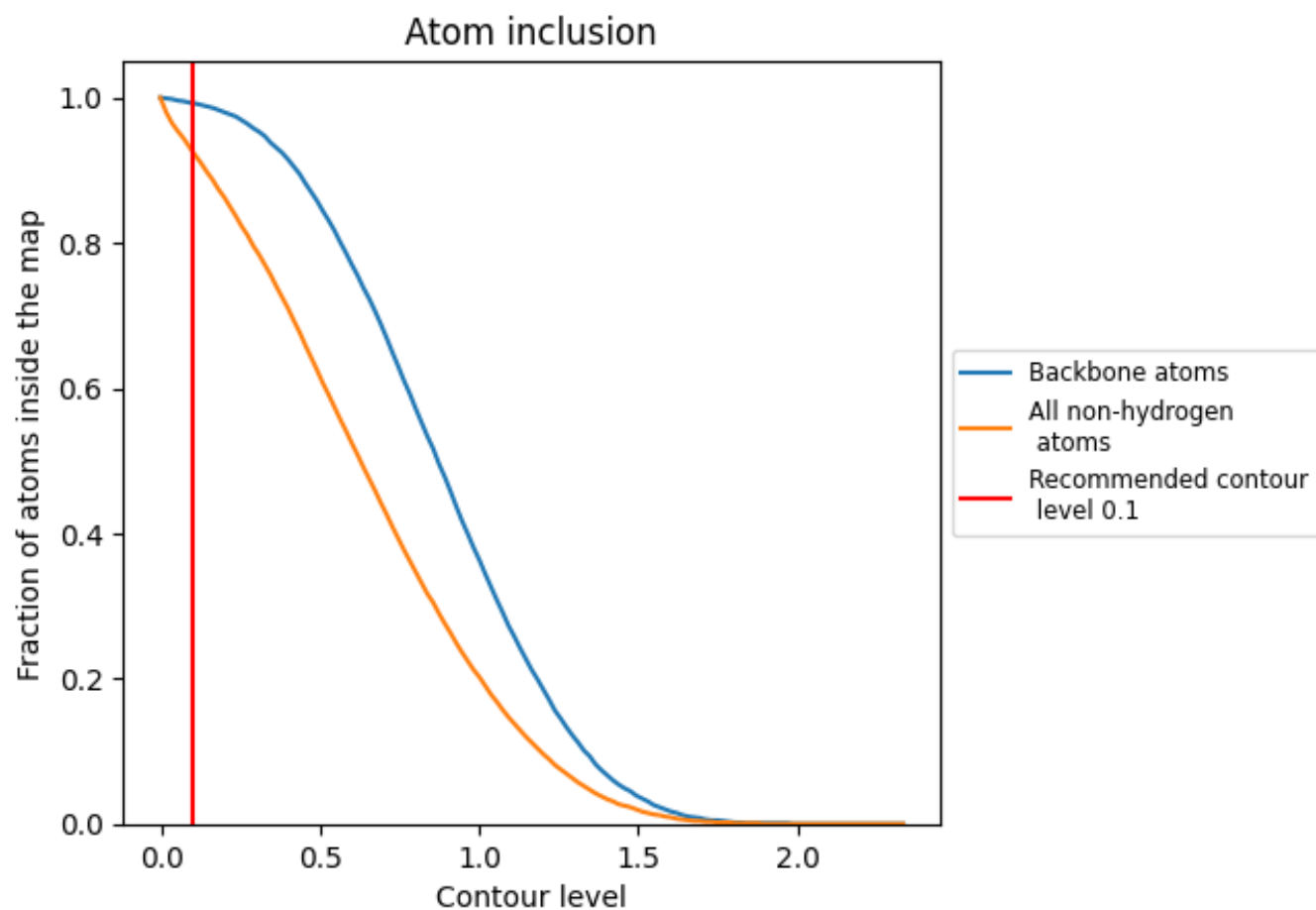
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 99% of all backbone atoms, 92% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9250	 0.2740
A	 0.9220	 0.2630
B	 0.9340	 0.2930
C	 0.9210	 0.2790
D	 0.7500	 0.1720
E	 0.8930	 0.1650
F	 0.9640	 0.2300
G	 0.9290	 0.2890
H	 0.9640	 0.2930
I	 0.8950	 0.2070
J	 0.9490	 0.2530
K	 0.9490	 0.1140
L	 1.0000	 0.1240
M	 0.9490	 0.0450
N	 0.9290	 0.1970
O	 0.9290	 0.1830
P	 0.8210	 -0.0690
Q	 0.9290	 0.1680
R	 0.8930	 0.1390
S	 1.0000	 0.3660
T	 0.8570	 0.1470
U	 1.0000	 0.2070
V	 0.9490	 0.1310
W	 1.0000	 0.1180
X	 0.9740	 0.2740
Y	 0.5360	 0.0540
Z	 0.9290	 0.0430
a	 0.9290	 0.2570
b	 1.0000	 0.2000
c	 0.9640	 0.0570
d	 0.9290	 0.2900
e	 0.8930	 0.2550
f	 0.9470	 0.3630
g	 0.9740	 0.1910
h	 0.9490	 0.2030



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Chain	Atom inclusion	Q-score
i	 0.8720	 0.1240
j	 0.9740	 0.1450