



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 12:34 PM UTC

PDB ID : 7SBT / pdb_00007sbt
EMDB ID : EMD-24988
Title : One RBD-up 2 of pre-fusion SARS-CoV-2 Gamma 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 : 4.40 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:

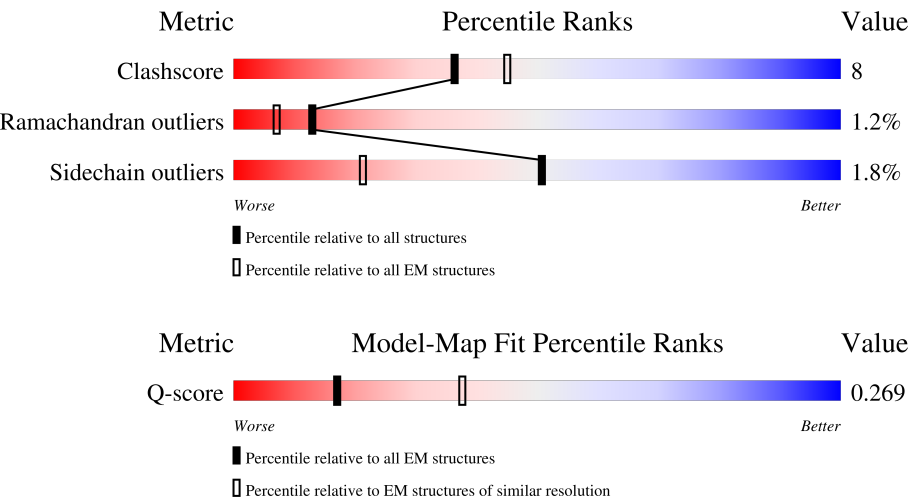
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data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

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

The reported resolution of this entry is 4.40 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	3132 (3.91 - 4.90)

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	74%9%16%
1	B	1310	76%6%16%
1	C	1310	76%8%16%
2	D	2	100%

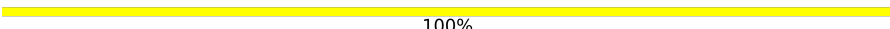
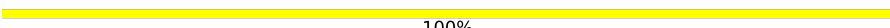
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Mol	Chain	Length	Quality of chain
2	E	2	100%
2	F	2	100%
2	G	2	100%
2	L	2	100%
2	M	2	50% 50%
2	N	2	50% 50%
2	O	2	50% 50%
2	P	2	100%
2	Q	2	100%
2	V	2	100%
2	W	2	100%
2	X	2	50% 50%
2	Y	2	100%
2	Z	2	100%
2	a	2	100%
2	b	2	100%
3	H	3	33% 67%
3	R	3	100%
3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

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Mol	Chain	Length	Quality of chain
4	d	3	 67%33%
4	e	3	 100%
4	f	3	 100%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	X	2	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 27151 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	1104	Total	C	N	O	S	0	0
			8627	5511	1431	1646	39		
1	B	1096	Total	C	N	O	S	0	0
			8582	5484	1426	1634	38		
1	C	1100	Total	C	N	O	S	0	0
			8609	5500	1430	1640	39		

There are 147 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	PHE	LEU	variant	UNP P0DTC2
A	20	ASN	THR	variant	UNP P0DTC2
A	26	SER	PRO	variant	UNP P0DTC2
A	138	TYR	ASP	variant	UNP P0DTC2
A	190	SER	ARG	variant	UNP P0DTC2
A	417	THR	LYS	variant	UNP P0DTC2
A	484	LYS	GLU	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	655	TYR	HIS	variant	UNP P0DTC2
A	1027	ILE	THR	variant	UNP P0DTC2
A	1176	PHE	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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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	18	PHE	LEU	variant	UNP P0DTC2
B	20	ASN	THR	variant	UNP P0DTC2
B	26	SER	PRO	variant	UNP P0DTC2
B	138	TYR	ASP	variant	UNP P0DTC2
B	190	SER	ARG	variant	UNP P0DTC2
B	417	THR	LYS	variant	UNP P0DTC2
B	484	LYS	GLU	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	1027	ILE	THR	variant	UNP P0DTC2
B	1176	PHE	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

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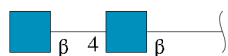
Chain	Residue	Modelled	Actual	Comment	Reference
B	1279	GLY	-	expression tag	UNP P0DTC2
B	1280	SER	-	expression tag	UNP P0DTC2
B	1281	ALA	-	expression tag	UNP P0DTC2
B	1282	TRP	-	expression tag	UNP P0DTC2
B	1283	SER	-	expression tag	UNP P0DTC2
B	1284	HIS	-	expression tag	UNP P0DTC2
B	1285	PRO	-	expression tag	UNP P0DTC2
B	1286	GLN	-	expression tag	UNP P0DTC2
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	18	PHE	LEU	variant	UNP P0DTC2
C	20	ASN	THR	variant	UNP P0DTC2
C	26	SER	PRO	variant	UNP P0DTC2
C	138	TYR	ASP	variant	UNP P0DTC2
C	190	SER	ARG	variant	UNP P0DTC2
C	417	THR	LYS	variant	UNP P0DTC2
C	484	LYS	GLU	variant	UNP P0DTC2
C	501	TYR	ASN	variant	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	655	TYR	HIS	variant	UNP P0DTC2

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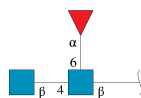
Chain	Residue	Modelled	Actual	Comment	Reference
C	1027	ILE	THR	variant	UNP P0DTC2
C	1176	PHE	VAL	variant	UNP P0DTC2
C	1274	LEU	-	expression tag	UNP P0DTC2
C	1275	GLU	-	expression tag	UNP P0DTC2
C	1276	SER	-	expression tag	UNP P0DTC2
C	1277	GLY	-	expression tag	UNP P0DTC2
C	1278	GLY	-	expression tag	UNP 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
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		
2	G	2	Total	C	N	O	0	0
			28	16	2	10		
2	L	2	Total	C	N	O	0	0
			28	16	2	10		
2	M	2	Total	C	N	O	0	0
			28	16	2	10		
2	N	2	Total	C	N	O	0	0
			28	16	2	10		
2	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	V	2	Total	C	N	O	0	0
			28	16	2	10		
2	W	2	Total	C	N	O	0	0
			28	16	2	10		
2	X	2	Total	C	N	O	0	0
			28	16	2	10		
2	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		

- 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	H	3	Total	C	N	O	0	0
			38	22	2	14		
3	R	3	Total	C	N	O	0	0
			38	22	2	14		
3	c	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	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		
4	K	3	Total	C	N	O	0	0
			39	22	2	15		
4	S	3	Total	C	N	O	0	0
			39	22	2	15		
4	T	3	Total	C	N	O	0	0
			39	22	2	15		
4	U	3	Total	C	N	O	0	0
			39	22	2	15		
4	d	3	Total	C	N	O	0	0
			39	22	2	15		
4	e	3	Total	C	N	O	0	0
			39	22	2	15		
4	f	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	

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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	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	
5	C	1	Total	C	N	O	0
			14	8	1	5	

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 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 L:  100%MAG1
MAG2

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

Chain M:  50% 50%MAG1
MAG2

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

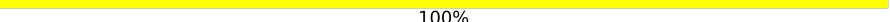
Chain N:  50% 50%MAG1
MAG2

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

Chain O:  50% 50%

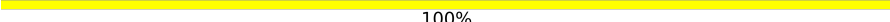
MAG1
MAG2

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

Chain P:  100%

MAG1
MAG2

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

Chain Q:  100%

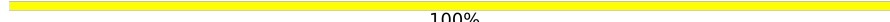
MAG1
MAG2

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

Chain V:  100%

MAG1
MAG2

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

Chain W:  100%

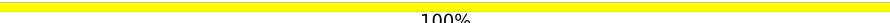
MAG1
MAG2

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

Chain X:  50% 50%

MAG1
MAG2

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

Chain Y:  100%

MAG1
MAG2

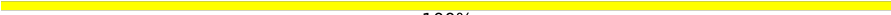
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z:  100%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 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  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 R:  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

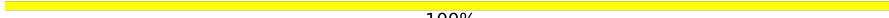
Chain c:  67% 33%MAG1
MAG2
FUC3

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

Chain I:  67% 33%

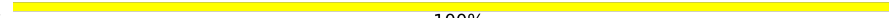
MAG1
MAG2
MAN3

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

Chain J:  100%

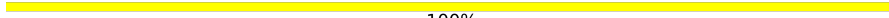
MAG1
MAG2
MAN3

- 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:  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 S:  100%

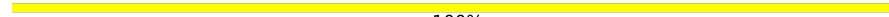
MAG1
MAG2
MAN3

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

Chain T:  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 U:  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 d:  67% 33%



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

Chain e:  100%



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

Chain f:  100%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	36346	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 ($e^-/\text{\AA}^2$)	54.72	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	2.175	Depositor
Minimum map value	-0.002	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.049	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: NAG, MAN, 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	0.70	1/8827 (0.0%)	1.26	21/12007 (0.2%)
1	B	0.69	0/8783	1.25	14/11949 (0.1%)
1	C	0.70	0/8810	1.25	12/11986 (0.1%)
All	All	0.70	1/26420 (0.0%)	1.25	47/35942 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	18	PHE	C-N	-7.98	1.22	1.33

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	164	ASN	N-CA-C	10.83	123.08	111.28
1	B	138	TYR	CA-C-N	8.13	128.14	119.76
1	B	138	TYR	C-N-CA	8.13	128.14	119.76
1	A	80	ASP	CA-CB-CG	6.81	119.41	112.60
1	B	80	ASP	CA-CB-CG	6.50	119.10	112.60
1	A	370	ASN	CA-CB-CG	6.39	118.99	112.60
1	B	458	LYS	N-CA-C	6.36	118.29	111.36
1	B	544	ASN	CA-CB-CG	6.22	118.82	112.60
1	A	833	PHE	N-CA-C	6.21	118.94	109.62
1	A	1107	ARG	N-CA-C	6.17	120.36	112.34
1	A	122	ASN	CA-CB-CG	5.97	118.57	112.60
1	B	635	VAL	N-CA-C	5.96	121.74	109.34
1	A	1110	TYR	CA-C-N	5.89	127.47	122.28
1	A	1110	TYR	C-N-CA	5.89	127.47	122.28
1	C	615	VAL	CA-C-O	-5.76	113.72	120.75
1	C	170	TYR	N-CA-C	5.74	117.40	107.93
1	A	342	PHE	O-C-N	5.66	129.92	122.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	619	GLU	CA-C-N	5.63	123.79	120.24
1	A	619	GLU	C-N-CA	5.63	123.79	120.24
1	C	76	THR	CA-C-N	5.52	127.94	120.65
1	C	76	THR	C-N-CA	5.52	127.94	120.65
1	B	1084	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	347	PHE	CA-CB-CG	5.50	119.30	113.80
1	B	394	ASN	OD1-CG-ND2	-5.49	117.11	122.60
1	C	73	THR	N-CA-C	5.48	117.52	109.24
1	B	979	ASP	CA-CB-CG	5.45	118.05	112.60
1	C	1127	ASP	CA-CB-CG	5.45	118.05	112.60
1	B	422	ASN	N-CA-C	5.44	116.96	109.29
1	C	106	PHE	CA-CB-CG	5.38	119.18	113.80
1	B	122	ASN	CA-CB-CG	5.36	117.96	112.60
1	C	122	ASN	CA-CB-CG	5.35	117.95	112.60
1	C	1064	HIS	CB-CG-CD2	-5.32	124.29	131.20
1	B	40	ASP	CA-CB-CG	5.31	117.91	112.60
1	A	913	GLN	OE1-CD-NE2	-5.28	117.32	122.60
1	A	333	THR	N-CA-C	5.28	117.11	111.36
1	A	178	ASP	CA-CB-CG	5.26	117.86	112.60
1	A	520	ALA	N-CA-CB	-5.22	105.91	111.29
1	A	145	TYR	N-CA-C	5.20	117.38	108.90
1	C	615	VAL	N-CA-CB	5.17	117.67	111.31
1	B	559	PHE	CA-CB-CG	5.14	118.94	113.80
1	A	1083	HIS	CB-CG-CD2	-5.12	124.54	131.20
1	A	1064	HIS	CB-CG-CD2	-5.11	124.56	131.20
1	C	565	PHE	CA-C-N	5.09	125.11	121.65
1	C	565	PHE	C-N-CA	5.09	125.11	121.65
1	A	995	ARG	N-CA-C	-5.04	105.44	111.03
1	A	23	GLN	N-CA-C	5.03	117.17	109.07
1	A	455	LEU	N-CA-CB	-5.02	104.22	110.90

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	8627	0	8386	169	0
1	B	8582	0	8346	157	0
1	C	8609	0	8368	123	0
2	D	28	0	25	2	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	L	28	0	25	0	0
2	M	28	0	25	3	0
2	N	28	0	25	0	0
2	O	28	0	25	0	0
2	P	28	0	25	0	0
2	Q	28	0	25	0	0
2	V	28	0	25	3	0
2	W	28	0	25	1	0
2	X	28	0	25	8	0
2	Y	28	0	25	0	0
2	Z	28	0	25	4	0
2	a	28	0	25	0	0
2	b	28	0	25	0	0
3	H	38	0	34	4	0
3	R	38	0	34	0	0
3	c	38	0	34	1	0
4	I	39	0	34	1	0
4	J	39	0	34	0	0
4	K	39	0	34	0	0
4	S	39	0	34	0	0
4	T	39	0	34	1	0
4	U	39	0	34	1	0
4	d	39	0	34	1	0
4	e	39	0	34	0	0
4	f	39	0	34	1	0
5	A	154	0	143	1	0
5	B	126	0	117	7	0
5	C	112	0	104	1	0
All	All	27151	0	26297	411	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 8.

All (411) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:497:PHE:CE1	1:A:507:PRO:HB3	1.34	1.57
1:B:1107:ARG:HD3	1:C:904:TYR:CE2	1.55	1.41
1:B:811:LYS:HB3	1:B:812:PRO:CD	1.56	1.31
1:C:578:ASP:OD1	1:C:579:PRO:HD2	1.37	1.24
1:A:497:PHE:HE1	1:A:507:PRO:CB	1.51	1.23
1:B:1107:ARG:CD	1:C:904:TYR:HE2	1.52	1.23
1:A:497:PHE:CE1	1:A:507:PRO:CB	2.21	1.22
1:B:57:PRO:HB3	1:B:273:ARG:NH1	1.56	1.18
1:A:834:ILE:HD12	2:Z:1:NAG:H81	1.26	1.17
1:A:377:PHE:CD2	1:A:434:ILE:HG12	1.78	1.17
1:B:520:ALA:CB	1:B:521:PRO:CD	2.23	1.15
1:B:811:LYS:HB3	1:B:812:PRO:HD3	1.29	1.14
1:B:813:SER:HB3	1:B:868:GLU:HG2	1.23	1.13
1:B:133:PHE:CD2	1:B:135:PHE:CZ	2.37	1.12
1:B:133:PHE:CG	1:B:135:PHE:CZ	2.37	1.11
1:C:339:GLY:O	1:C:343:ASN:HB2	1.49	1.11
1:A:457:ARG:HH22	1:A:461:LEU:HB3	1.15	1.10
1:B:520:ALA:HB1	1:B:521:PRO:CD	1.79	1.10
1:B:24:LEU:HB3	1:B:25:PRO:CD	1.80	1.09
1:B:133:PHE:CD2	1:B:135:PHE:HZ	1.70	1.08
1:A:457:ARG:NH2	1:A:461:LEU:HB3	1.68	1.08
1:B:520:ALA:CB	1:B:521:PRO:HD3	1.83	1.07
1:B:133:PHE:HB2	1:B:135:PHE:CE2	1.91	1.06
1:C:140:PHE:CZ	1:C:158:ARG:HD2	1.91	1.05
1:C:577:ARG:HH21	1:C:582:LEU:HD22	1.10	1.04
1:B:24:LEU:CB	1:B:25:PRO:HD3	1.88	1.04
1:B:1107:ARG:CD	1:C:904:TYR:CE2	2.31	1.03
1:B:28:TYR:HD1	1:B:61:ASN:OD1	1.42	1.02
1:B:610:VAL:HG23	1:B:636:TYR:CE2	1.93	1.02
1:B:811:LYS:HB3	1:B:812:PRO:HD2	1.41	1.02
1:A:447:GLY:HA2	1:A:497:PHE:O	1.58	1.01
1:B:520:ALA:HB1	1:B:521:PRO:HD3	1.05	1.01
1:B:133:PHE:CB	1:B:135:PHE:CZ	2.43	1.00
1:B:808:ASP:HB3	1:B:809:PRO:HD2	1.40	0.98
1:B:136:CYS:O	1:B:139:PRO:HD3	1.61	0.98
1:B:1107:ARG:HD3	1:C:904:TYR:CD2	1.99	0.97
1:A:319:ARG:HD2	1:B:740:MET:HE3	1.47	0.96
1:A:377:PHE:HD2	1:A:434:ILE:HG12	1.16	0.96
1:B:811:LYS:CB	1:B:812:PRO:CD	2.42	0.95
1:A:377:PHE:CD2	1:A:434:ILE:CG1	2.48	0.95
1:A:319:ARG:HD2	1:B:740:MET:CE	1.96	0.95
1:B:133:PHE:HB2	1:B:135:PHE:CZ	2.03	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:740:MET:HG2	1:C:857:GLY:HA3	1.45	0.93
1:B:24:LEU:HB3	1:B:25:PRO:HD3	0.94	0.93
1:B:718:PHE:CG	1:B:1067:TYR:CE2	2.56	0.92
1:C:740:MET:CG	1:C:857:GLY:HA3	1.98	0.92
1:B:133:PHE:CG	1:B:135:PHE:CE1	2.58	0.91
1:A:904:TYR:CD1	1:C:1107:ARG:HD3	2.07	0.90
1:A:18:PHE:CE2	1:A:138:TYR:HE2	1.90	0.89
1:A:1016:ALA:HA	1:A:1019:ARG:NH1	1.87	0.89
1:B:1014:ARG:HG2	1:B:1014:ARG:HH11	1.37	0.89
1:B:147:LYS:HB3	5:B:1402:NAG:H81	1.53	0.88
1:B:147:LYS:HB3	5:B:1402:NAG:C8	2.02	0.88
1:B:813:SER:CB	1:B:868:GLU:HG2	2.03	0.88
1:B:520:ALA:HB3	1:B:521:PRO:CD	2.05	0.87
2:D:1:NAG:H62	2:D:2:NAG:HN2	1.40	0.86
1:B:520:ALA:HB3	1:B:521:PRO:HD2	1.55	0.86
1:A:406:GLU:HG2	1:A:418:ILE:HG13	1.59	0.85
1:C:953:ASN:O	1:C:957:GLN:HG3	1.76	0.85
1:A:896:ILE:HD11	1:A:904:TYR:CE2	2.10	0.85
1:A:1016:ALA:HA	1:A:1019:ARG:HH12	1.39	0.85
1:B:23:GLN:HA	1:B:80:ASP:HB3	1.58	0.85
1:A:331:ASN:OD1	1:A:580:GLN:CB	2.24	0.84
1:B:610:VAL:CG2	1:B:636:TYR:CE2	2.61	0.84
1:B:900:MET:HG3	1:B:904:TYR:CE1	2.12	0.84
1:B:807:PRO:HB3	1:B:814:LYS:O	1.78	0.83
1:A:453:TYR:CD2	1:A:493:GLN:O	2.32	0.83
1:A:18:PHE:CZ	1:A:258:TRP:NE1	2.47	0.83
1:C:577:ARG:NH2	1:C:582:LEU:HD22	1.93	0.83
1:A:24:LEU:HD13	1:A:24:LEU:O	1.80	0.82
1:B:811:LYS:CB	1:B:812:PRO:HD3	2.06	0.82
1:C:339:GLY:O	1:C:343:ASN:CB	2.28	0.81
1:A:904:TYR:CE1	1:C:1107:ARG:HD3	2.15	0.81
1:B:357:ARG:HH21	1:C:230:PRO:HG3	1.44	0.81
1:A:453:TYR:CE2	1:A:493:GLN:O	2.34	0.81
1:C:896:ILE:HD11	1:C:904:TYR:CE1	2.15	0.80
1:B:808:ASP:HB3	1:B:809:PRO:CD	2.12	0.80
1:A:331:ASN:OD1	1:A:580:GLN:HB3	1.82	0.80
1:A:25:PRO:HD2	1:A:80:ASP:OD2	1.82	0.80
1:A:148:ASN:HB2	5:A:1403:NAG:H82	1.64	0.80
1:A:904:TYR:CE1	1:C:1107:ARG:CD	2.64	0.79
1:A:377:PHE:CE2	1:A:434:ILE:HD11	2.18	0.79
1:B:900:MET:HG3	1:B:904:TYR:HE1	1.45	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:ARG:CD	1:B:740:MET:CE	2.61	0.78
1:C:741:TYR:OH	1:C:962:LEU:HD12	1.83	0.78
1:B:458:LYS:HZ2	2:X:2:NAG:C6	1.96	0.78
1:A:457:ARG:HD2	1:A:459:SER:O	1.83	0.78
1:C:1014:ARG:HB2	1:C:1014:ARG:NH1	1.97	0.78
1:A:831:ALA:HB3	1:A:833:PHE:CE1	2.19	0.78
2:M:1:NAG:H61	2:M:2:NAG:C7	2.14	0.77
1:B:896:ILE:HD11	1:B:904:TYR:CE1	2.20	0.77
1:A:18:PHE:CE2	1:A:138:TYR:CE2	2.71	0.76
1:A:18:PHE:HE2	1:A:138:TYR:HE2	1.32	0.76
1:C:740:MET:HG2	1:C:857:GLY:CA	2.15	0.76
1:A:831:ALA:HB3	1:A:833:PHE:HE1	1.51	0.76
1:B:718:PHE:CG	1:B:1067:TYR:HE2	2.00	0.76
1:B:608:VAL:HG23	1:B:636:TYR:OH	1.86	0.76
2:D:1:NAG:H62	2:D:2:NAG:N2	2.00	0.75
1:A:24:LEU:H	1:A:25:PRO:CD	1.99	0.75
1:C:577:ARG:HH21	1:C:582:LEU:CD2	1.92	0.75
1:B:1107:ARG:HD3	1:C:904:TYR:HE2	0.95	0.75
1:A:834:ILE:HD12	2:Z:1:NAG:C8	2.13	0.75
1:B:1107:ARG:CG	1:C:904:TYR:HE2	2.00	0.75
1:B:458:LYS:NZ	2:X:2:NAG:H62	2.02	0.75
1:C:741:TYR:CE2	1:C:1004:LEU:HG	2.22	0.75
1:C:900:MET:O	1:C:904:TYR:CD1	2.39	0.74
1:C:900:MET:O	1:C:904:TYR:HD1	1.71	0.74
1:A:498:GLN:HB3	1:A:500:THR:HG22	1.69	0.74
1:C:271:GLN:OE1	1:C:273:ARG:NH1	2.20	0.74
1:A:332:ILE:HD12	1:A:332:ILE:H	1.53	0.73
1:B:1039:ARG:NH2	1:C:1031:GLU:OE2	2.21	0.73
1:C:741:TYR:OH	1:C:962:LEU:CD1	2.36	0.73
1:C:905:ARG:HH11	1:C:1036:GLN:HB2	1.54	0.73
1:B:140:PHE:CZ	1:B:158:ARG:HD2	2.23	0.73
1:B:28:TYR:CD1	1:B:61:ASN:OD1	2.35	0.73
1:A:145:TYR:O	1:A:147:LYS:HG2	1.88	0.73
1:A:131:CYS:HA	1:A:165:ASN:O	1.90	0.72
1:A:904:TYR:HE1	1:C:1107:ARG:CD	2.03	0.72
1:B:458:LYS:NZ	2:X:2:NAG:C6	2.52	0.72
1:B:458:LYS:HZ2	2:X:2:NAG:H61	1.54	0.72
1:B:133:PHE:CB	1:B:135:PHE:CE2	2.69	0.72
1:B:900:MET:O	1:B:904:TYR:CD1	2.43	0.72
1:A:377:PHE:CE2	1:A:434:ILE:HG12	2.26	0.71
1:A:1101:HIS:CE1	4:I:1:NAG:H3	2.25	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:497:PHE:HE1	1:A:507:PRO:CA	2.03	0.71
1:A:900:MET:HG3	1:A:904:TYR:HE2	1.54	0.71
1:C:905:ARG:NH1	1:C:1036:GLN:HB2	2.04	0.71
1:A:904:TYR:CE1	1:C:1107:ARG:NE	2.59	0.71
1:B:813:SER:O	1:B:815:ARG:N	2.23	0.70
1:A:452:LEU:HA	1:A:494:SER:HA	1.72	0.70
1:A:332:ILE:HD12	1:A:332:ILE:N	2.06	0.70
1:A:457:ARG:HH22	1:A:461:LEU:CB	1.99	0.70
1:A:377:PHE:CE2	1:A:434:ILE:CG1	2.75	0.70
1:B:57:PRO:HB3	1:B:273:ARG:HH11	1.55	0.70
1:A:24:LEU:HD13	1:A:24:LEU:C	2.16	0.69
1:C:171:VAL:HG21	2:V:2:NAG:H82	1.73	0.69
1:C:741:TYR:CD2	1:C:1004:LEU:HG	2.27	0.69
1:B:1039:ARG:NE	1:C:1031:GLU:OE2	2.25	0.69
1:A:452:LEU:HD23	1:A:494:SER:HB3	1.75	0.69
1:C:896:ILE:HD11	1:C:904:TYR:HE1	1.57	0.69
1:A:331:ASN:OD1	1:A:580:GLN:CG	2.42	0.68
1:A:377:PHE:HE1	1:A:384:PRO:HB3	1.58	0.68
1:B:133:PHE:CB	1:B:135:PHE:CE1	2.75	0.68
1:C:131:CYS:HA	1:C:165:ASN:O	1.95	0.67
1:A:740:MET:SD	1:C:592:PHE:CE1	2.87	0.67
1:B:357:ARG:NH2	1:C:230:PRO:HD3	2.09	0.67
1:A:904:TYR:CD1	1:C:1107:ARG:CD	2.77	0.67
1:B:147:LYS:O	1:B:149:ASN:N	2.21	0.67
1:C:741:TYR:CZ	1:C:962:LEU:CD1	2.78	0.67
1:C:900:MET:HG3	1:C:904:TYR:HE1	1.60	0.66
1:A:773:GLU:OE2	1:A:1019:ARG:NH2	2.28	0.66
1:B:131:CYS:HA	1:B:165:ASN:O	1.95	0.66
1:C:741:TYR:CZ	1:C:962:LEU:HD11	2.30	0.66
1:A:900:MET:HG3	1:A:904:TYR:CE2	2.31	0.66
1:A:146:HIS:CE1	1:A:150:LYS:HG3	2.31	0.66
1:A:23:GLN:HE22	1:A:82:PRO:HB3	1.61	0.65
1:A:834:ILE:HG23	2:Z:1:NAG:O7	1.95	0.65
1:B:357:ARG:HB2	1:B:396:TYR:CE1	2.31	0.65
1:A:377:PHE:CE2	1:A:434:ILE:CD1	2.80	0.64
1:A:319:ARG:HG2	1:B:740:MET:HE1	1.79	0.64
1:C:326:ILE:HD12	1:C:539:VAL:HG21	1.79	0.64
1:A:448:ASN:H	1:A:497:PHE:HB2	1.63	0.63
1:B:610:VAL:HG23	1:B:636:TYR:CZ	2.33	0.63
1:A:28:TYR:HB3	1:A:61:ASN:OD1	1.97	0.63
1:B:1014:ARG:HH11	1:B:1014:ARG:CG	2.11	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:443:SER:HA	1:A:497:PHE:HB3	1.80	0.63
1:B:996:LEU:O	1:B:1000:ARG:HG3	1.99	0.63
1:C:741:TYR:CE1	1:C:962:LEU:HG	2.33	0.62
1:A:497:PHE:HE1	1:A:507:PRO:HB3	0.87	0.62
1:A:332:ILE:H	1:A:332:ILE:CD1	2.12	0.62
1:C:331:ASN:OD1	1:C:331:ASN:O	2.16	0.62
1:B:79:PHE:HE2	1:B:138:TYR:OH	1.83	0.62
1:C:132:GLU:OE1	1:C:165:ASN:ND2	2.33	0.61
1:A:406:GLU:HG2	1:A:418:ILE:CG1	2.31	0.61
1:B:357:ARG:HH22	1:C:230:PRO:HD3	1.66	0.61
1:B:415:THR:HG23	1:B:415:THR:O	2.01	0.61
1:B:357:ARG:HB2	1:B:396:TYR:HE1	1.66	0.61
1:A:148:ASN:O	1:A:149:ASN:OD1	2.19	0.60
1:B:458:LYS:HZ1	2:X:2:NAG:H62	1.66	0.60
1:A:145:TYR:HB2	1:A:152:TRP:CH2	2.36	0.60
1:A:146:HIS:HE1	1:A:150:LYS:HG3	1.64	0.60
1:B:61:ASN:ND2	5:B:1401:NAG:O7	2.35	0.60
1:B:147:LYS:HB3	5:B:1402:NAG:C7	2.30	0.60
1:B:422:ASN:HD21	1:B:454:ARG:HB3	1.65	0.60
1:C:905:ARG:HD2	1:C:1049:LEU:O	2.01	0.60
1:C:900:MET:HG3	1:C:904:TYR:CE1	2.36	0.60
1:A:497:PHE:CZ	1:A:507:PRO:HB3	2.22	0.60
1:B:57:PRO:CB	1:B:273:ARG:NH1	2.50	0.60
1:A:145:TYR:HB3	1:A:152:TRP:CZ3	2.37	0.60
1:A:18:PHE:HE2	1:A:138:TYR:CE2	2.14	0.60
1:B:900:MET:O	1:B:904:TYR:HD1	1.83	0.60
1:C:741:TYR:CD2	1:C:1004:LEU:CG	2.85	0.59
1:C:578:ASP:OD1	1:C:579:PRO:CD	2.31	0.59
1:C:740:MET:HG3	1:C:857:GLY:HA3	1.82	0.59
1:A:497:PHE:CD1	1:A:507:PRO:CG	2.86	0.59
1:B:147:LYS:CB	5:B:1402:NAG:O7	2.50	0.59
1:A:983:ARG:NH2	1:C:430:THR:HG21	2.18	0.59
1:A:497:PHE:HD1	1:A:507:PRO:HD3	1.66	0.59
1:A:905:ARG:NH1	1:A:1036:GLN:HB2	2.17	0.59
1:C:330:PRO:HA	1:C:579:PRO:HB2	1.84	0.58
1:A:365:TYR:HE2	1:A:388:ASN:N	2.01	0.58
1:A:904:TYR:HD1	1:C:1107:ARG:HD3	1.66	0.58
1:A:24:LEU:N	1:A:25:PRO:CD	2.66	0.58
1:A:377:PHE:HE2	1:A:434:ILE:HD11	1.63	0.58
2:X:1:NAG:O4	2:X:2:NAG:O7	2.20	0.58
1:A:452:LEU:CD2	1:A:494:SER:HB3	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:497:PHE:CE1	1:A:507:PRO:CG	2.86	0.58
1:B:718:PHE:CD1	1:B:1067:TYR:CE2	2.91	0.58
1:B:896:ILE:HD11	1:B:904:TYR:HE1	1.66	0.58
1:A:352:ALA:O	1:A:466:ARG:NH2	2.36	0.58
1:B:811:LYS:HD2	1:B:811:LYS:N	2.18	0.58
1:C:14:GLN:O	1:C:158:ARG:HG2	2.04	0.57
1:A:365:TYR:HE2	1:A:388:ASN:H	1.53	0.57
1:B:79:PHE:CE2	1:B:138:TYR:OH	2.56	0.57
1:C:154:GLU:HB3	2:V:1:NAG:H81	1.86	0.57
1:A:23:GLN:NE2	1:A:82:PRO:HB3	2.18	0.57
1:A:24:LEU:H	1:A:25:PRO:HD2	1.67	0.57
1:B:147:LYS:HB2	5:B:1402:NAG:O7	2.04	0.57
1:B:742:ILE:HG23	1:B:1000:ARG:HD2	1.86	0.57
1:A:23:GLN:HG3	1:A:80:ASP:OD2	2.05	0.57
1:C:578:ASP:HB3	1:C:581:THR:O	2.05	0.57
1:C:132:GLU:OE1	1:C:165:ASN:HB2	2.04	0.57
1:A:896:ILE:HD11	1:A:904:TYR:HE2	1.63	0.56
1:C:1098:ASN:OD1	1:C:1101:HIS:N	2.26	0.56
1:A:319:ARG:CD	1:B:740:MET:HE2	2.34	0.56
1:C:1014:ARG:HB2	1:C:1014:ARG:HH11	1.70	0.56
1:A:904:TYR:CD1	1:C:1107:ARG:NE	2.74	0.56
1:B:357:ARG:NH2	1:C:230:PRO:HG3	2.16	0.56
1:B:357:ARG:HH21	1:C:230:PRO:CG	2.18	0.56
1:B:421:TYR:CE2	1:B:456:PHE:HA	2.40	0.56
1:B:1014:ARG:HG2	1:B:1014:ARG:NH1	2.15	0.56
1:A:453:TYR:CZ	1:A:493:GLN:HG3	2.40	0.56
1:A:905:ARG:HH11	1:A:1036:GLN:HB2	1.70	0.56
1:C:578:ASP:CG	1:C:579:PRO:HD2	2.23	0.56
1:B:133:PHE:CD1	1:B:135:PHE:CE1	2.93	0.56
1:A:133:PHE:CB	1:A:135:PHE:CZ	2.89	0.55
1:B:1107:ARG:NE	1:C:904:TYR:CE2	2.74	0.55
1:B:57:PRO:HB3	1:B:273:ARG:HH12	1.63	0.55
1:C:140:PHE:HE2	1:C:158:ARG:HH11	1.55	0.55
1:B:457:ARG:C	1:B:473:TYR:CE1	2.84	0.55
1:B:133:PHE:HB3	1:B:135:PHE:CE1	2.41	0.55
1:A:406:GLU:HG3	1:A:409:GLN:NE2	2.22	0.55
1:C:339:GLY:O	1:C:343:ASN:CA	2.55	0.55
1:A:145:TYR:HB2	1:A:152:TRP:CZ2	2.42	0.54
1:C:154:GLU:HB3	2:V:1:NAG:C8	2.37	0.54
1:A:834:ILE:CG2	2:Z:1:NAG:O7	2.55	0.54
1:B:409:GLN:CD	1:B:416:GLY:HA3	2.32	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:497:PHE:CD1	1:A:507:PRO:HD3	2.42	0.54
1:B:520:ALA:CB	1:B:521:PRO:HD2	2.13	0.54
1:C:886:TRP:CE3	1:C:905:ARG:NH1	2.76	0.54
1:C:737:ASP:O	1:C:740:MET:HB3	2.08	0.54
1:B:458:LYS:NZ	2:X:2:NAG:H61	2.18	0.54
1:A:145:TYR:CB	1:A:152:TRP:CZ3	2.91	0.54
1:A:447:GLY:CA	1:A:497:PHE:O	2.45	0.54
1:A:497:PHE:CD1	1:A:507:PRO:HG3	2.42	0.54
1:C:1014:ARG:HH11	1:C:1014:ARG:CB	2.21	0.53
1:A:377:PHE:CD2	1:A:434:ILE:HG13	2.42	0.53
1:A:1094:VAL:CG2	1:B:904:TYR:OH	2.57	0.53
4:T:1:NAG:O4	4:T:2:NAG:O7	2.26	0.53
1:A:23:GLN:HB2	1:A:80:ASP:HB3	1.91	0.53
1:B:74:ASN:OD1	4:U:1:NAG:O5	2.23	0.53
1:B:418:ILE:O	1:B:422:ASN:HB2	2.08	0.53
1:B:718:PHE:CB	1:B:1067:TYR:CE2	2.92	0.53
1:B:718:PHE:CD2	1:B:1067:TYR:HE2	2.27	0.52
1:B:1039:ARG:CZ	1:C:1031:GLU:OE2	2.56	0.52
1:C:74:ASN:ND2	4:f:1:NAG:C7	2.71	0.52
1:C:454:ARG:HD3	1:C:457:ARG:HG2	1.92	0.52
1:A:748:GLU:H	1:A:748:GLU:CD	2.17	0.52
1:B:742:ILE:O	1:B:742:ILE:HG22	2.08	0.52
1:B:357:ARG:HD2	1:B:396:TYR:OH	2.10	0.52
1:B:79:PHE:HE2	1:B:138:TYR:HH	1.48	0.52
1:C:158:ARG:HH21	1:C:254:SER:C	2.18	0.52
1:A:319:ARG:HE	1:B:740:MET:HE2	1.75	0.52
1:B:357:ARG:CB	1:B:396:TYR:CE1	2.93	0.52
1:B:742:ILE:HG12	1:B:1000:ARG:HB3	1.91	0.51
1:B:422:ASN:ND2	1:B:454:ARG:HB3	2.25	0.51
1:C:454:ARG:HH22	1:C:469:SER:HB3	1.76	0.51
2:X:1:NAG:H61	2:X:2:NAG:C7	2.41	0.51
1:A:331:ASN:OD1	1:A:580:GLN:CA	2.59	0.51
1:C:355:ARG:HG3	1:C:398:ASP:OD1	2.10	0.51
3:H:1:NAG:H62	3:H:2:NAG:O7	2.11	0.51
1:A:377:PHE:HE2	1:A:434:ILE:CD1	2.22	0.51
1:A:145:TYR:HB3	1:A:152:TRP:CE3	2.46	0.51
1:A:404:GLY:C	1:A:406:GLU:H	2.20	0.50
1:A:904:TYR:HE1	1:C:1107:ARG:NE	2.05	0.50
1:C:74:ASN:N	1:C:74:ASN:OD1	2.43	0.50
3:H:1:NAG:H61	3:H:2:NAG:C7	2.41	0.50
1:A:133:PHE:HB2	1:A:135:PHE:CZ	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:ARG:NE	1:B:740:MET:HE2	2.27	0.50
1:C:175:PHE:CE1	1:C:177:MET:HE2	2.47	0.50
1:C:886:TRP:HE3	1:C:905:ARG:NH1	2.10	0.50
1:A:404:GLY:HA3	1:A:508:TYR:CE2	2.48	0.49
1:A:904:TYR:CD1	1:C:1107:ARG:CZ	2.95	0.49
1:B:295:PRO:HG2	1:B:636:TYR:CE1	2.46	0.49
1:B:1039:ARG:HE	1:C:1031:GLU:CD	2.19	0.49
1:A:831:ALA:CB	1:A:833:PHE:HE1	2.24	0.49
1:A:332:ILE:N	1:A:332:ILE:CD1	2.73	0.49
1:B:418:ILE:O	1:B:422:ASN:C	2.56	0.49
1:A:365:TYR:OH	1:A:385:THR:O	2.30	0.48
1:C:164:ASN:ND2	1:C:165:ASN:CG	2.72	0.48
1:A:403:ARG:HG2	1:A:504:GLY:O	2.14	0.48
1:B:905:ARG:HD2	1:B:1049:LEU:O	2.13	0.48
1:A:493:GLN:OE1	1:A:493:GLN:HA	2.14	0.48
1:B:357:ARG:NH2	1:C:230:PRO:CD	2.77	0.48
1:B:357:ARG:NH2	1:C:230:PRO:CG	2.77	0.48
1:B:421:TYR:HE2	1:B:457:ARG:H	1.62	0.47
1:B:147:LYS:CB	5:B:1402:NAG:C7	2.92	0.47
1:B:457:ARG:C	1:B:473:TYR:HE1	2.23	0.47
1:C:1014:ARG:HB2	1:C:1014:ARG:CZ	2.44	0.47
1:A:326:ILE:HG23	1:A:541:PHE:HA	1.97	0.47
1:A:133:PHE:HB2	1:A:135:PHE:CE2	2.50	0.47
1:A:444:LYS:O	1:A:497:PHE:O	2.32	0.47
1:A:983:ARG:NH2	1:C:381:GLY:O	2.48	0.46
1:B:808:ASP:CB	1:B:809:PRO:CD	2.84	0.46
1:C:326:ILE:HD11	1:C:534:VAL:HG22	1.97	0.46
1:C:762:GLN:HA	1:C:765:ARG:HE	1.81	0.46
1:A:319:ARG:CG	1:B:740:MET:CE	2.93	0.46
1:A:453:TYR:CZ	1:A:493:GLN:CG	2.99	0.46
1:A:983:ARG:NH2	1:C:430:THR:CG2	2.78	0.46
1:B:455:LEU:HG	1:B:456:PHE:CD2	2.51	0.46
3:H:1:NAG:C6	3:H:2:NAG:C7	2.93	0.46
1:B:357:ARG:CD	1:B:396:TYR:OH	2.63	0.46
1:C:592:PHE:O	1:C:592:PHE:CG	2.68	0.46
1:C:593:GLY:HA3	1:C:615:VAL:CG2	2.46	0.45
1:A:331:ASN:OD1	1:A:580:GLN:HA	2.16	0.45
1:A:133:PHE:CG	1:A:135:PHE:CZ	3.04	0.45
1:A:365:TYR:CD2	1:A:387:LEU:HB2	2.51	0.45
1:B:761:THR:O	1:B:765:ARG:HG3	2.17	0.45
1:C:592:PHE:O	1:C:592:PHE:CD2	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:392:PHE:CD2	1:C:517:LEU:HD21	2.52	0.45
1:C:748:GLU:CD	1:C:748:GLU:H	2.23	0.45
1:B:357:ARG:CG	1:B:396:TYR:CE1	3.00	0.45
1:A:24:LEU:C	1:A:24:LEU:CD1	2.86	0.45
1:B:1014:ARG:CG	1:B:1014:ARG:NH1	2.72	0.45
1:C:886:TRP:CE3	1:C:905:ARG:CZ	3.00	0.45
1:A:369:TYR:HD1	1:A:377:PHE:CD1	2.35	0.45
1:A:979:ASP:O	1:A:983:ARG:HD3	2.17	0.45
1:A:1094:VAL:HB	1:B:904:TYR:OH	2.17	0.44
1:C:213:VAL:O	1:C:213:VAL:HG23	2.16	0.44
1:B:357:ARG:HG3	1:B:396:TYR:CE1	2.53	0.44
1:B:748:GLU:CD	1:B:748:GLU:H	2.25	0.44
1:C:1014:ARG:NH1	1:C:1014:ARG:CB	2.73	0.44
1:A:452:LEU:HD23	1:A:494:SER:CB	2.47	0.44
1:B:1107:ARG:CD	1:C:904:TYR:CD2	2.85	0.44
1:A:796:ASP:OD2	1:C:709:ASN:N	2.47	0.44
2:M:1:NAG:O4	2:M:2:NAG:O7	2.34	0.44
1:A:1107:ARG:HG2	1:B:904:TYR:HE2	1.83	0.44
1:C:520:ALA:HB1	1:C:521:PRO:HD2	1.99	0.44
1:A:497:PHE:CD1	1:A:507:PRO:CD	3.00	0.44
1:B:158:ARG:NH2	1:B:254:SER:HA	2.33	0.44
1:A:498:GLN:HE21	1:A:500:THR:CG2	2.31	0.44
1:A:1011:GLN:OE1	1:A:1014:ARG:NH1	2.51	0.44
3:c:1:NAG:H62	3:c:2:NAG:H82	1.99	0.44
1:A:146:HIS:O	1:A:146:HIS:CD2	2.70	0.44
1:A:740:MET:SD	1:C:592:PHE:CD1	3.11	0.43
1:B:24:LEU:HD13	1:B:24:LEU:C	2.43	0.43
1:A:23:GLN:CB	1:A:80:ASP:HB3	2.47	0.43
1:A:145:TYR:CB	1:A:152:TRP:CE3	3.01	0.43
1:B:422:ASN:HD21	1:B:454:ARG:CB	2.31	0.43
1:C:1101:HIS:CE1	4:d:1:NAG:H3	2.54	0.43
1:A:904:TYR:HE1	1:C:1107:ARG:CG	2.32	0.43
1:B:138:TYR:O	1:B:138:TYR:CG	2.70	0.43
1:A:497:PHE:HD1	1:A:507:PRO:CD	2.32	0.43
1:B:104:TRP:HB2	1:B:106:PHE:CE1	2.53	0.43
1:A:404:GLY:CA	1:A:508:TYR:CD2	3.01	0.43
1:A:831:ALA:HB3	1:A:833:PHE:CD1	2.52	0.43
1:B:357:ARG:HG3	1:B:396:TYR:CZ	2.54	0.43
1:B:900:MET:CG	1:B:904:TYR:HE1	2.22	0.43
1:C:593:GLY:HA3	1:C:615:VAL:HG21	2.00	0.43
1:A:905:ARG:HD2	1:A:1049:LEU:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:ASN:HB3	1:B:423:TYR:H	1.55	0.42
1:B:1039:ARG:CZ	1:B:1042:PHE:CD1	3.02	0.42
1:C:339:GLY:O	1:C:343:ASN:N	2.52	0.42
1:B:718:PHE:CD2	1:B:1067:TYR:CE2	3.04	0.42
1:B:813:SER:OG	1:B:815:ARG:NH1	2.52	0.42
1:B:742:ILE:HG23	1:B:1000:ARG:CD	2.50	0.42
1:C:164:ASN:ND2	1:C:165:ASN:ND2	2.67	0.42
1:A:24:LEU:N	1:A:25:PRO:HD3	2.35	0.42
1:A:904:TYR:HE1	1:C:1107:ARG:HG2	1.85	0.42
1:C:164:ASN:ND2	2:W:1:NAG:O5	2.46	0.42
1:C:1090:PRO:HD3	1:C:1095:PHE:CE2	2.55	0.42
1:A:136:CYS:O	1:A:139:PRO:HD3	2.20	0.41
1:B:610:VAL:CG2	1:B:636:TYR:CD2	3.03	0.41
1:A:453:TYR:HB3	1:A:495:TYR:CE2	2.55	0.41
1:B:64:TRP:CD1	1:B:266:TYR:CE2	3.09	0.41
1:B:1039:ARG:NH1	1:B:1042:PHE:CE1	2.88	0.41
1:A:1031:GLU:CD	1:C:1039:ARG:HE	2.28	0.41
1:A:1158:ASN:O	1:A:1162:PRO:HD2	2.21	0.41
1:A:453:TYR:OH	1:A:493:GLN:HG3	2.19	0.41
1:A:868:GLU:H	1:A:868:GLU:CD	2.29	0.41
1:A:145:TYR:HB2	1:A:152:TRP:CZ3	2.55	0.41
1:A:457:ARG:HD3	1:A:459:SER:H	1.86	0.41
1:A:136:CYS:C	1:A:138:TYR:N	2.78	0.41
1:A:643:PHE:CZ	1:A:655:TYR:CD1	3.09	0.41
1:B:158:ARG:HH21	1:B:254:SER:HB2	1.85	0.41
1:B:815:ARG:HG2	1:B:871:ALA:HB2	2.02	0.41
1:C:422:ASN:HD22	1:C:454:ARG:HB3	1.86	0.41
1:A:319:ARG:HG2	1:B:740:MET:CE	2.46	0.41
1:A:339:GLY:O	1:A:343:ASN:HB2	2.21	0.41
1:B:133:PHE:HB3	1:B:135:PHE:CD1	2.55	0.41
1:A:904:TYR:HD1	1:C:1107:ARG:CD	2.29	0.40
1:C:710:ASN:HB3	5:C:1407:NAG:H82	2.03	0.40
1:A:365:TYR:CE2	1:A:388:ASN:N	2.87	0.40
1:B:149:ASN:HB2	1:B:151:SER:OG	2.21	0.40
1:C:896:ILE:CD1	1:C:904:TYR:CE1	2.97	0.40
2:M:1:NAG:H61	2:M:2:NAG:C8	2.51	0.40
1:C:21:ARG:CZ	1:C:138:TYR:HB2	2.50	0.40
1:A:900:MET:O	1:A:904:TYR:CD2	2.74	0.40
3:H:1:NAG:C6	3:H:2:NAG:O7	2.69	0.40
1:C:602:THR:O	1:C:604:THR:N	2.54	0.40
1:C:886:TRP:HZ3	1:C:905:ARG:HD3	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1094/1310 (84%)	1004 (92%)	74 (7%)	16 (2%)	8	38
1	B	1086/1310 (83%)	988 (91%)	86 (8%)	12 (1%)	11	45
1	C	1090/1310 (83%)	995 (91%)	85 (8%)	10 (1%)	14	49
All	All	3270/3930 (83%)	2987 (91%)	245 (8%)	38 (1%)	13	42

All (38) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	920	GLN
1	B	148	ASN
1	B	520	ALA
1	B	635	VAL
1	B	812	PRO
1	B	814	LYS
1	C	603	ASN
1	A	74	ASN
1	A	150	LYS
1	A	744	GLY
1	B	24	LEU
1	B	811	LYS
1	C	32	PHE
1	A	59	PHE
1	A	79	PHE
1	A	137	ASN
1	A	709	ASN
1	B	423	TYR
1	B	709	ASN
1	C	24	LEU

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Mol	Chain	Res	Type
1	C	940	SER
1	C	1041	ASP
1	A	405	ASP
1	A	985	ASP
1	A	1041	ASP
1	B	79	PHE
1	B	59	PHE
1	A	283	GLY
1	A	332	ILE
1	B	25	PRO
1	C	518	LEU
1	C	709	ASN
1	A	834	ILE
1	A	367	VAL
1	C	579	PRO
1	C	159	VAL
1	C	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	962/1135 (85%)	939 (98%)	23 (2%)	43	63
1	B	958/1135 (84%)	948 (99%)	10 (1%)	68	76
1	C	961/1135 (85%)	941 (98%)	20 (2%)	47	65
All	All	2881/3405 (85%)	2828 (98%)	53 (2%)	51	67

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	16	VAL
1	A	46	SER
1	A	74	ASN
1	A	80	ASP

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Mol	Chain	Res	Type
1	A	122	ASN
1	A	186	PHE
1	A	282	ASN
1	A	284	THR
1	A	332	ILE
1	A	335	LEU
1	A	351	TYR
1	A	367	VAL
1	A	368	LEU
1	A	370	ASN
1	A	406	GLU
1	A	461	LEU
1	A	514	SER
1	A	515	PHE
1	A	603	ASN
1	A	787	GLN
1	A	803	SER
1	A	850	ILE
1	B	24	LEU
1	B	77	LYS
1	B	122	ASN
1	B	186	PHE
1	B	376	THR
1	B	441	LEU
1	B	457	ARG
1	B	544	ASN
1	B	803	SER
1	B	965	GLN
1	C	18	PHE
1	C	73	THR
1	C	74	ASN
1	C	122	ASN
1	C	185	ASN
1	C	186	PHE
1	C	214	ARG
1	C	332	ILE
1	C	422	ASN
1	C	461	LEU
1	C	488	CYS
1	C	523	THR
1	C	618	THR
1	C	622	VAL

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Mol	Chain	Res	Type
1	C	624	ILE
1	C	760	CYS
1	C	901	GLN
1	C	965	GLN
1	C	1072	GLU
1	C	1127	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	23	GLN
1	A	121	ASN
1	A	125	ASN
1	A	146	HIS
1	A	183	GLN
1	A	271	GLN
1	A	354	ASN
1	A	360	ASN
1	A	409	GLN
1	A	498	GLN
1	A	675	GLN
1	A	804	GLN
1	A	919	ASN
1	A	935	GLN
1	A	949	GLN
1	A	960	ASN
1	A	1054	GLN
1	A	1088	HIS
1	A	1135	ASN
1	B	23	GLN
1	B	121	ASN
1	B	164	ASN
1	B	354	ASN
1	B	422	ASN
1	B	439	ASN
1	B	450	ASN
1	B	856	ASN
1	B	872	GLN
1	B	907	ASN
1	B	919	ASN
1	B	953	ASN

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Mol	Chain	Res	Type
1	B	955	ASN
1	B	1002	GLN
1	B	1011	GLN
1	B	1119	ASN
1	C	137	ASN
1	C	164	ASN
1	C	360	ASN
1	C	580	GLN
1	C	613	GLN
1	C	992	GLN
1	C	1002	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 ⓘ

70 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	1.20	1 (7%)	17,19,21	1.17	2 (11%)
2	NAG	D	2	2	14,14,15	1.12	1 (7%)	17,19,21	0.83	1 (5%)
2	NAG	E	1	2,1	14,14,15	1.20	1 (7%)	17,19,21	1.07	2 (11%)
2	NAG	E	2	2	14,14,15	1.28	2 (14%)	17,19,21	0.90	0
2	NAG	F	1	2,1	14,14,15	1.25	1 (7%)	17,19,21	1.03	2 (11%)
2	NAG	F	2	2	14,14,15	1.26	2 (14%)	17,19,21	0.90	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	G	1	2,1	14,14,15	1.21	1 (7%)	17,19,21	0.91	0
2	NAG	G	2	2	14,14,15	1.21	1 (7%)	17,19,21	0.67	0
3	NAG	H	1	1,3	14,14,15	1.29	1 (7%)	17,19,21	0.97	2 (11%)
3	NAG	H	2	3	14,14,15	1.26	2 (14%)	17,19,21	1.20	1 (5%)
3	FUC	H	3	3	10,10,11	1.81	2 (20%)	14,14,16	1.21	2 (14%)
4	NAG	I	1	4,1	14,14,15	1.33	2 (14%)	17,19,21	1.20	3 (17%)
4	NAG	I	2	4	14,14,15	1.51	4 (28%)	17,19,21	0.97	1 (5%)
4	MAN	I	3	4	11,11,12	1.45	2 (18%)	15,15,17	0.81	0
4	NAG	J	1	4,1	14,14,15	1.14	1 (7%)	17,19,21	0.62	0
4	NAG	J	2	4	14,14,15	1.21	2 (14%)	17,19,21	1.08	1 (5%)
4	MAN	J	3	4	11,11,12	1.41	2 (18%)	15,15,17	0.93	0
4	NAG	K	1	4,1	14,14,15	1.32	2 (14%)	17,19,21	2.19	3 (17%)
4	NAG	K	2	4	14,14,15	1.74	4 (28%)	17,19,21	1.73	2 (11%)
4	MAN	K	3	4	11,11,12	1.42	2 (18%)	15,15,17	1.07	1 (6%)
2	NAG	L	1	2,1	14,14,15	1.35	3 (21%)	17,19,21	0.82	1 (5%)
2	NAG	L	2	2	14,14,15	1.29	2 (14%)	17,19,21	0.84	1 (5%)
2	NAG	M	1	2,1	14,14,15	0.24	0	17,19,21	0.52	0
2	NAG	M	2	2	14,14,15	1.14	1 (7%)	17,19,21	1.27	1 (5%)
2	NAG	N	1	2,1	14,14,15	0.27	0	17,19,21	0.45	0
2	NAG	N	2	2	14,14,15	1.17	1 (7%)	17,19,21	0.95	1 (5%)
2	NAG	O	1	2,1	14,14,15	0.29	0	17,19,21	0.46	0
2	NAG	O	2	2	14,14,15	1.15	1 (7%)	17,19,21	1.18	1 (5%)
2	NAG	P	1	2,1	14,14,15	1.08	1 (7%)	17,19,21	0.98	1 (5%)
2	NAG	P	2	2	14,14,15	1.19	1 (7%)	17,19,21	0.97	1 (5%)
2	NAG	Q	1	2,1	14,14,15	1.16	1 (7%)	17,19,21	1.04	1 (5%)
2	NAG	Q	2	2	14,14,15	1.47	2 (14%)	17,19,21	0.68	0
3	NAG	R	1	1,3	14,14,15	0.26	0	17,19,21	0.91	1 (5%)
3	NAG	R	2	3	14,14,15	1.06	1 (7%)	17,19,21	0.94	1 (5%)
3	FUC	R	3	3	10,10,11	1.49	1 (10%)	14,14,16	1.50	3 (21%)
4	NAG	S	1	4,1	14,14,15	1.13	1 (7%)	17,19,21	1.08	2 (11%)
4	NAG	S	2	4	14,14,15	1.25	2 (14%)	17,19,21	0.70	0
4	MAN	S	3	4	11,11,12	1.26	2 (18%)	15,15,17	0.92	0
4	NAG	T	1	4,1	14,14,15	1.54	2 (14%)	17,19,21	1.00	1 (5%)
4	NAG	T	2	4	14,14,15	1.27	1 (7%)	17,19,21	1.58	2 (11%)
4	MAN	T	3	4	11,11,12	1.55	2 (18%)	15,15,17	0.92	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	U	1	4,1	14,14,15	0.23	0	17,19,21	0.50	0
4	NAG	U	2	4	14,14,15	1.17	2 (14%)	17,19,21	1.28	3 (17%)
4	MAN	U	3	4	11,11,12	1.44	2 (18%)	15,15,17	0.93	1 (6%)
2	NAG	V	1	2,1	14,14,15	1.34	3 (21%)	17,19,21	0.83	1 (5%)
2	NAG	V	2	2	14,14,15	1.28	2 (14%)	17,19,21	0.84	1 (5%)
2	NAG	W	1	2,1	14,14,15	0.32	0	17,19,21	0.58	0
2	NAG	W	2	2	14,14,15	1.01	1 (7%)	17,19,21	0.81	0
2	NAG	X	1	2,1	14,14,15	0.30	0	17,19,21	0.56	0
2	NAG	X	2	2	14,14,15	1.03	1 (7%)	17,19,21	1.22	1 (5%)
2	NAG	Y	1	2,1	14,14,15	1.32	1 (7%)	17,19,21	1.47	3 (17%)
2	NAG	Y	2	2	14,14,15	1.28	1 (7%)	17,19,21	0.99	1 (5%)
2	NAG	Z	1	2,1	14,14,15	0.21	0	17,19,21	0.45	0
2	NAG	Z	2	2	14,14,15	1.07	1 (7%)	17,19,21	0.78	0
2	NAG	a	1	2,1	14,14,15	1.21	2 (14%)	17,19,21	0.66	0
2	NAG	a	2	2	14,14,15	1.34	2 (14%)	17,19,21	1.05	1 (5%)
2	NAG	b	1	2,1	14,14,15	0.97	1 (7%)	17,19,21	0.95	1 (5%)
2	NAG	b	2	2	14,14,15	1.16	1 (7%)	17,19,21	0.90	1 (5%)
3	NAG	c	1	1,3	14,14,15	0.20	0	17,19,21	0.84	1 (5%)
3	NAG	c	2	3	14,14,15	0.29	0	17,19,21	0.54	0
3	FUC	c	3	3	10,10,11	1.58	2 (20%)	14,14,16	1.30	2 (14%)
4	NAG	d	1	4,1	14,14,15	1.15	2 (14%)	17,19,21	1.08	1 (5%)
4	NAG	d	2	4	14,14,15	1.41	3 (21%)	17,19,21	0.92	1 (5%)
4	MAN	d	3	4	11,11,12	1.47	2 (18%)	15,15,17	0.75	0
4	NAG	e	1	4,1	14,14,15	1.12	1 (7%)	17,19,21	0.95	0
4	NAG	e	2	4	14,14,15	1.31	2 (14%)	17,19,21	0.68	0
4	MAN	e	3	4	11,11,12	1.57	2 (18%)	15,15,17	0.99	1 (6%)
4	NAG	f	1	4,1	14,14,15	0.24	0	17,19,21	0.51	0
4	NAG	f	2	4	14,14,15	1.21	2 (14%)	17,19,21	0.99	1 (5%)
4	MAN	f	3	4	11,11,12	1.42	2 (18%)	15,15,17	0.90	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	D	1	2,1	-	1/6/23/26	0/1/1/1

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

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

All (99) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	NAG	O5-C1	-4.41	1.36	1.43
4	T	1	NAG	O5-C5	3.91	1.51	1.43
4	T	3	MAN	O5-C5	3.49	1.50	1.43
4	e	3	MAN	O5-C5	3.47	1.50	1.43
4	I	2	NAG	O5-C5	3.25	1.49	1.43
2	N	2	NAG	O5-C5	3.22	1.49	1.43
2	F	2	NAG	O5-C5	3.14	1.49	1.43
2	M	2	NAG	O5-C5	3.11	1.49	1.43
4	K	2	NAG	O5-C1	3.11	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	3	FUC	O5-C5	3.10	1.49	1.43
4	I	1	NAG	O5-C5	3.10	1.49	1.43
3	c	3	FUC	C1-C2	3.04	1.59	1.52
3	H	3	FUC	O5-C1	3.04	1.48	1.43
2	a	2	NAG	O5-C5	3.03	1.49	1.43
2	Q	2	NAG	O5-C1	3.00	1.48	1.43
2	L	2	NAG	O5-C5	2.99	1.49	1.43
2	V	2	NAG	O5-C5	2.98	1.49	1.43
2	P	2	NAG	O5-C5	2.97	1.49	1.43
2	Q	2	NAG	O5-C5	2.97	1.49	1.43
4	d	3	MAN	O5-C5	2.96	1.49	1.43
2	F	1	NAG	O5-C5	2.94	1.49	1.43
4	K	2	NAG	O5-C5	2.89	1.49	1.43
2	E	2	NAG	O5-C5	2.87	1.49	1.43
4	J	3	MAN	O5-C5	2.85	1.49	1.43
2	O	2	NAG	O5-C5	2.85	1.49	1.43
2	Y	1	NAG	O5-C5	2.85	1.49	1.43
4	U	3	MAN	O5-C5	2.84	1.49	1.43
4	I	3	MAN	O5-C5	2.84	1.49	1.43
2	G	2	NAG	O5-C5	2.83	1.49	1.43
2	X	2	NAG	O5-C5	2.82	1.48	1.43
2	Z	2	NAG	O5-C5	2.80	1.48	1.43
3	c	3	FUC	O5-C5	2.79	1.49	1.43
3	H	2	NAG	O5-C5	2.78	1.48	1.43
2	L	1	NAG	O5-C5	2.77	1.48	1.43
4	d	2	NAG	O5-C5	2.77	1.48	1.43
4	K	3	MAN	O5-C5	2.76	1.48	1.43
2	Q	1	NAG	O5-C5	2.75	1.48	1.43
3	R	2	NAG	O5-C5	2.75	1.48	1.43
3	H	1	NAG	C1-C2	2.75	1.56	1.52
2	V	1	NAG	O5-C5	2.75	1.48	1.43
4	f	3	MAN	O5-C5	2.74	1.48	1.43
4	e	2	NAG	O5-C5	2.71	1.48	1.43
4	K	1	NAG	O5-C5	2.70	1.48	1.43
2	W	2	NAG	O5-C5	2.69	1.48	1.43
2	Y	2	NAG	O5-C5	2.66	1.48	1.43
2	E	1	NAG	O5-C5	2.63	1.48	1.43
2	D	2	NAG	O5-C5	2.62	1.48	1.43
4	e	3	MAN	O5-C1	2.61	1.48	1.43
2	a	1	NAG	O5-C5	2.61	1.48	1.43
4	K	2	NAG	O4-C4	2.60	1.49	1.43
4	U	2	NAG	O5-C5	2.59	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	R	3	FUC	O5-C5	2.57	1.48	1.43
4	U	3	MAN	O5-C1	2.56	1.48	1.43
4	K	2	NAG	C1-C2	2.56	1.55	1.52
4	J	2	NAG	O5-C5	2.53	1.48	1.43
4	S	3	MAN	O5-C5	2.50	1.48	1.43
4	d	2	NAG	O5-C1	2.50	1.47	1.43
4	T	2	NAG	O5-C5	2.50	1.48	1.43
4	T	3	MAN	O5-C1	2.45	1.47	1.43
4	f	2	NAG	O5-C5	2.45	1.48	1.43
4	f	2	NAG	O4-C4	2.44	1.49	1.43
4	S	2	NAG	O4-C4	2.43	1.49	1.43
4	S	2	NAG	O5-C5	2.42	1.48	1.43
2	b	2	NAG	O5-C5	2.41	1.48	1.43
4	S	1	NAG	O5-C5	2.40	1.48	1.43
4	e	1	NAG	O5-C5	2.40	1.48	1.43
2	P	1	NAG	O5-C5	2.39	1.48	1.43
4	T	1	NAG	C1-C2	2.38	1.55	1.52
4	I	3	MAN	O5-C1	2.37	1.47	1.43
4	K	3	MAN	O5-C1	2.37	1.47	1.43
2	a	2	NAG	O5-C1	2.32	1.47	1.43
3	H	2	NAG	C1-C2	2.30	1.55	1.52
4	e	2	NAG	O4-C4	2.29	1.48	1.43
4	J	1	NAG	O5-C5	2.29	1.47	1.43
4	d	1	NAG	O5-C1	2.29	1.47	1.43
4	I	2	NAG	C1-C2	2.29	1.55	1.52
4	J	2	NAG	O4-C4	2.29	1.48	1.43
4	d	3	MAN	O5-C1	2.29	1.47	1.43
4	I	2	NAG	O4-C4	2.26	1.48	1.43
4	J	3	MAN	O5-C1	2.24	1.47	1.43
4	d	1	NAG	O4-C4	2.22	1.48	1.43
2	L	1	NAG	O4-C4	2.19	1.48	1.43
2	V	1	NAG	O4-C4	2.18	1.48	1.43
2	E	2	NAG	O5-C1	2.17	1.47	1.43
4	I	2	NAG	O5-C1	2.16	1.47	1.43
4	f	3	MAN	O5-C1	2.16	1.47	1.43
2	V	1	NAG	C1-C2	2.13	1.55	1.52
2	L	1	NAG	C1-C2	2.13	1.55	1.52
4	U	2	NAG	O4-C4	2.12	1.48	1.43
2	b	1	NAG	O5-C5	2.11	1.47	1.43
4	S	3	MAN	O5-C1	2.07	1.47	1.43
2	G	1	NAG	O4-C4	2.06	1.48	1.43
4	I	1	NAG	O5-C1	2.06	1.47	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	a	1	NAG	O4-C4	2.04	1.48	1.43
2	V	2	NAG	O5-C1	2.04	1.47	1.43
4	K	1	NAG	O5-C1	2.04	1.47	1.43
2	F	2	NAG	O5-C1	2.03	1.47	1.43
2	L	2	NAG	O5-C1	2.03	1.47	1.43
4	d	2	NAG	C1-C2	2.02	1.55	1.52

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	K	1	NAG	C1-O5-C5	7.37	122.06	112.19
4	K	2	NAG	C4-C3-C2	5.32	118.82	111.02
2	M	2	NAG	C1-O5-C5	4.60	118.35	112.19
2	O	2	NAG	C1-O5-C5	4.24	117.87	112.19
4	T	2	NAG	C1-O5-C5	4.17	117.78	112.19
2	X	2	NAG	C1-O5-C5	4.12	117.71	112.19
3	H	2	NAG	C1-O5-C5	3.99	117.53	112.19
2	Y	1	NAG	C3-C4-C5	3.64	116.83	110.23
4	T	2	NAG	C4-C3-C2	-3.59	105.75	111.02
4	J	2	NAG	C1-O5-C5	3.36	116.68	112.19
4	K	2	NAG	O5-C5-C6	3.30	114.08	107.66
3	R	3	FUC	C1-O5-C5	3.28	120.70	112.97
4	K	1	NAG	C1-C2-N2	3.24	115.55	110.43
4	U	2	NAG	C2-N2-C7	-3.22	118.59	122.90
4	f	2	NAG	C1-O5-C5	3.20	116.48	112.19
4	d	1	NAG	C1-O5-C5	3.19	116.47	112.19
3	R	1	NAG	C1-O5-C5	3.17	116.43	112.19
3	R	2	NAG	C1-O5-C5	3.15	116.40	112.19
2	Y	1	NAG	C1-O5-C5	3.13	116.38	112.19
3	R	3	FUC	C1-C2-C3	3.04	114.08	109.64
2	a	2	NAG	C1-O5-C5	3.00	116.20	112.19
2	D	1	NAG	C1-O5-C5	2.97	116.17	112.19
3	c	3	FUC	C1-O5-C5	2.92	119.85	112.97
2	P	1	NAG	C4-C3-C2	-2.86	106.82	111.02
2	N	2	NAG	C1-O5-C5	2.75	115.87	112.19
3	c	1	NAG	C1-O5-C5	2.75	115.87	112.19
2	V	2	NAG	C1-O5-C5	2.73	115.84	112.19
2	L	2	NAG	C1-O5-C5	2.71	115.81	112.19
4	e	3	MAN	C1-O5-C5	2.68	115.77	112.19
4	K	1	NAG	C4-C3-C2	-2.62	107.17	111.02
4	I	1	NAG	C1-O5-C5	2.62	115.69	112.19
4	K	3	MAN	C1-O5-C5	2.57	115.62	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	b	1	NAG	C3-C4-C5	2.55	114.86	110.23
2	Y	2	NAG	C1-O5-C5	2.55	115.61	112.19
4	S	1	NAG	C3-C4-C5	2.55	114.85	110.23
3	c	3	FUC	C1-C2-C3	2.51	113.30	109.64
3	H	3	FUC	C1-C2-C3	2.50	113.28	109.64
2	P	2	NAG	C4-C3-C2	-2.49	107.37	111.02
2	D	2	NAG	C1-O5-C5	2.48	115.51	112.19
4	U	2	NAG	C1-C2-N2	2.48	114.34	110.43
2	E	1	NAG	C3-C4-C5	2.47	114.71	110.23
2	b	2	NAG	C3-C4-C5	2.37	114.53	110.23
3	H	1	NAG	O5-C5-C6	2.32	112.17	107.66
4	d	2	NAG	C4-C3-C2	-2.29	107.66	111.02
4	U	3	MAN	C1-O5-C5	2.27	115.23	112.19
2	Y	1	NAG	O5-C1-C2	-2.25	107.82	111.29
2	F	1	NAG	C1-O5-C5	2.24	115.19	112.19
4	S	1	NAG	O4-C4-C3	-2.14	105.32	110.38
3	R	3	FUC	C3-C4-C5	-2.13	106.58	109.81
3	H	1	NAG	C1-O5-C5	2.13	115.04	112.19
2	E	1	NAG	O5-C1-C2	-2.13	108.00	111.29
4	T	1	NAG	C4-C3-C2	-2.12	107.91	111.02
2	V	1	NAG	O5-C1-C2	-2.10	108.05	111.29
4	I	2	NAG	O5-C1-C2	-2.10	108.05	111.29
4	U	2	NAG	C1-O5-C5	2.09	114.99	112.19
2	L	1	NAG	O5-C1-C2	-2.08	108.07	111.29
2	D	1	NAG	C3-C4-C5	2.06	113.97	110.23
3	H	3	FUC	C1-O5-C5	2.03	117.77	112.97
2	F	1	NAG	O4-C4-C3	-2.03	105.59	110.38
2	Q	1	NAG	C1-O5-C5	-2.01	109.49	112.19
4	I	1	NAG	O5-C1-C2	-2.01	108.18	111.29
4	I	1	NAG	O4-C4-C3	-2.01	105.65	110.38

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	R	1	NAG	O5-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6
3	H	1	NAG	C4-C5-C6-O6
3	c	1	NAG	O5-C5-C6-O6
4	U	2	NAG	C8-C7-N2-C2
2	Y	1	NAG	C4-C5-C6-O6
3	R	1	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	K	2	NAG	C8-C7-N2-C2
3	c	1	NAG	C4-C5-C6-O6
4	K	2	NAG	C4-C5-C6-O6
4	U	2	NAG	O7-C7-N2-C2
4	K	2	NAG	O5-C5-C6-O6
4	U	1	NAG	O5-C5-C6-O6
4	f	1	NAG	O5-C5-C6-O6
4	S	3	MAN	O5-C5-C6-O6
4	d	3	MAN	O5-C5-C6-O6
4	U	1	NAG	C4-C5-C6-O6
4	f	1	NAG	C4-C5-C6-O6
2	Y	1	NAG	O5-C5-C6-O6
4	K	1	NAG	C8-C7-N2-C2
4	K	2	NAG	O7-C7-N2-C2
4	I	3	MAN	O5-C5-C6-O6
4	K	1	NAG	O7-C7-N2-C2
2	F	2	NAG	C1-C2-N2-C7
4	I	2	NAG	C1-C2-N2-C7
2	D	1	NAG	C3-C2-N2-C7
4	S	3	MAN	C4-C5-C6-O6

All (3) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	U	3	MAN	C1-C2-C3-C4-C5-O5
4	T	3	MAN	C1-C2-C3-C4-C5-O5
4	e	3	MAN	C1-C2-C3-C4-C5-O5

20 monomers are involved in 31 short contacts:

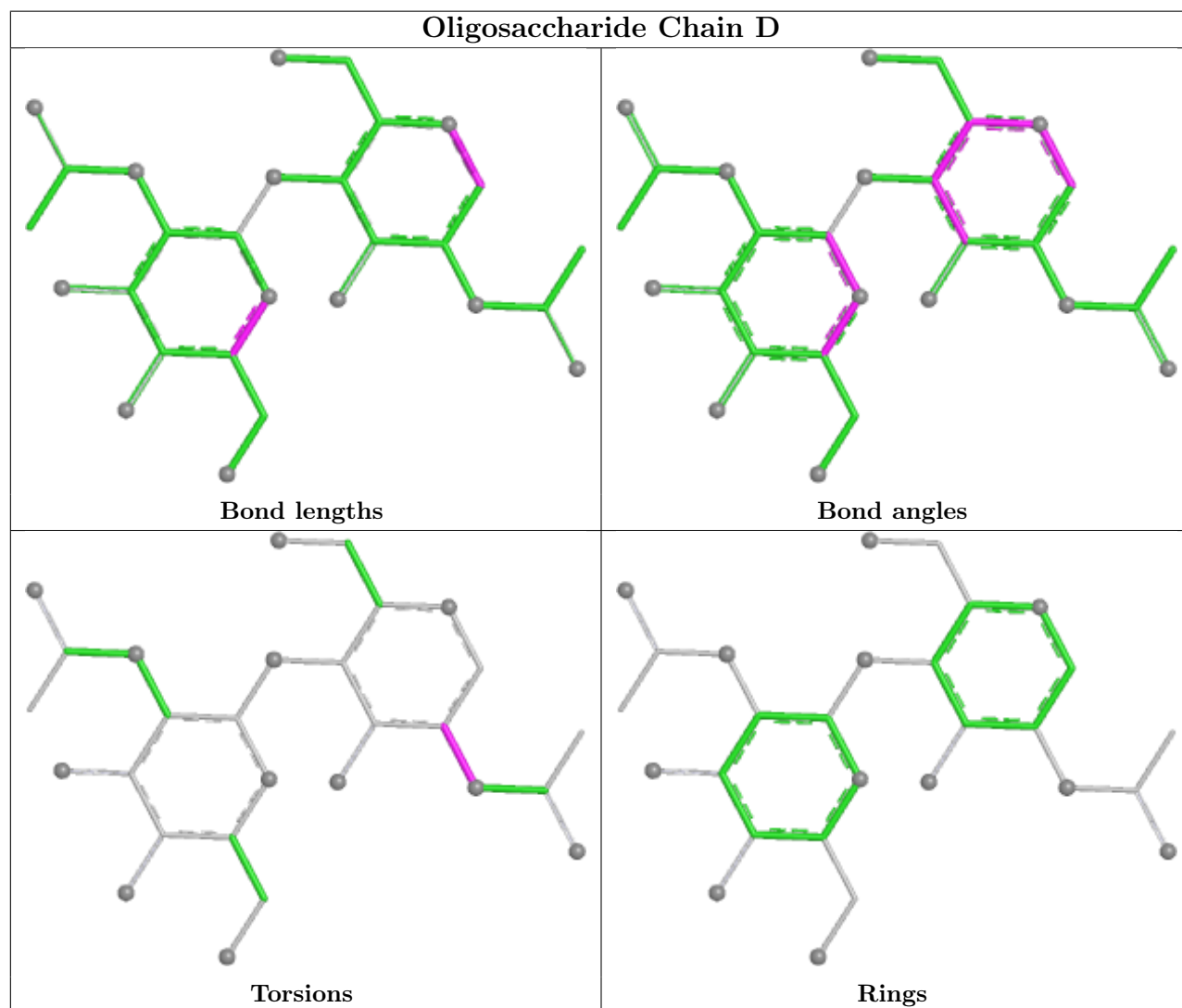
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	W	1	NAG	1	0
2	Z	1	NAG	4	0
4	f	1	NAG	1	0
2	V	1	NAG	2	0
2	X	2	NAG	8	0
4	T	1	NAG	1	0
4	T	2	NAG	1	0
2	D	2	NAG	2	0
2	D	1	NAG	2	0
2	M	2	NAG	3	0
2	V	2	NAG	1	0

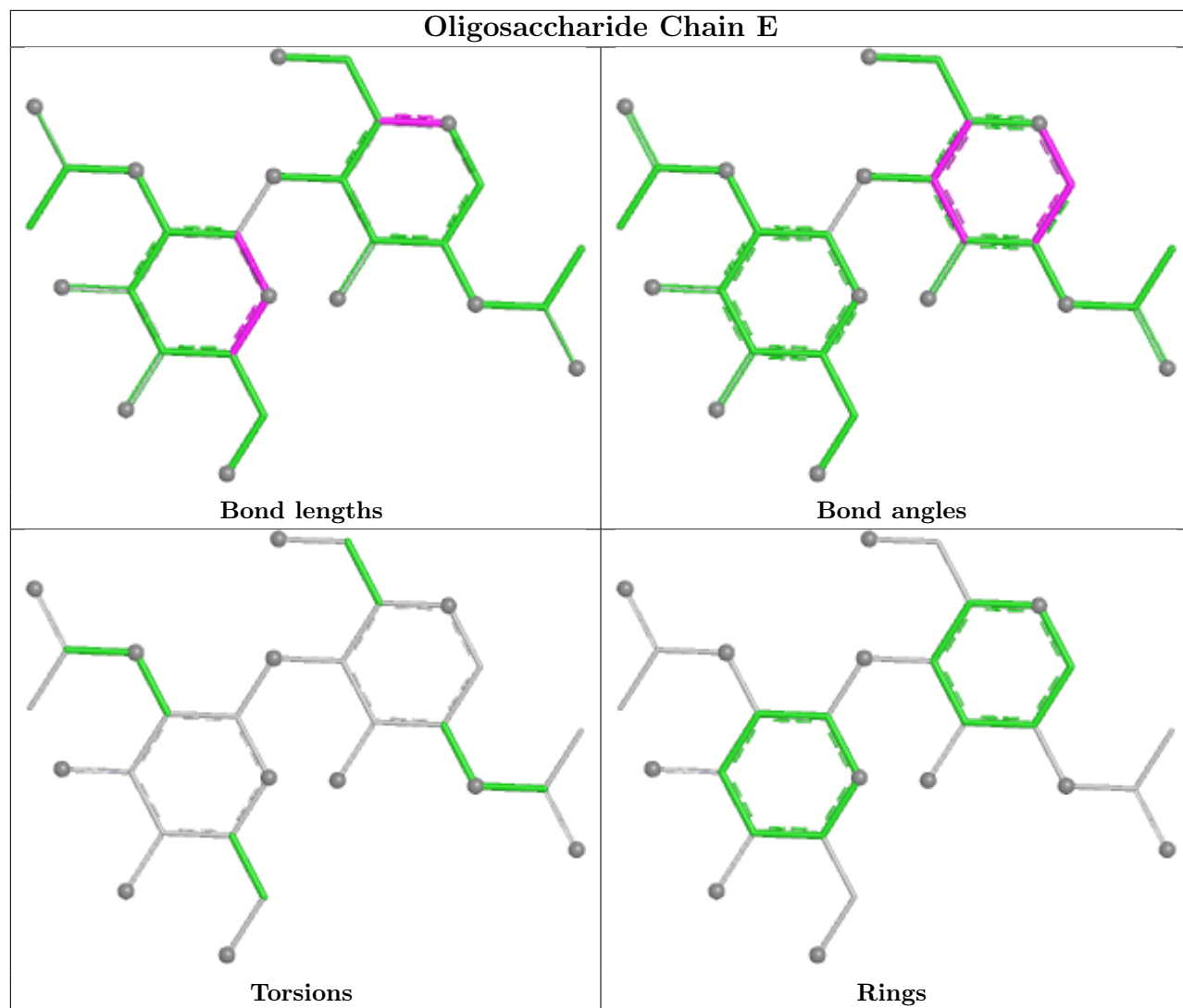
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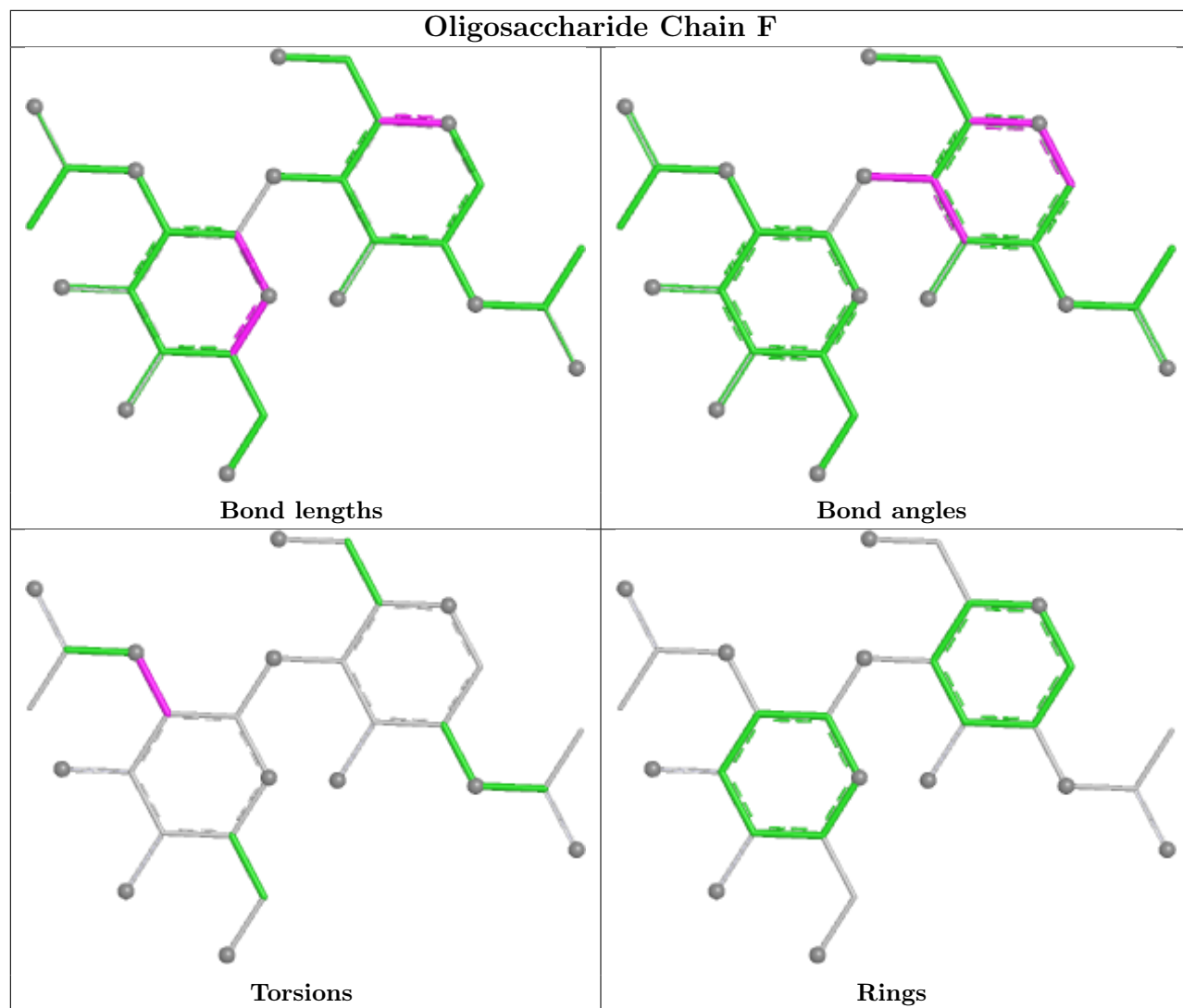
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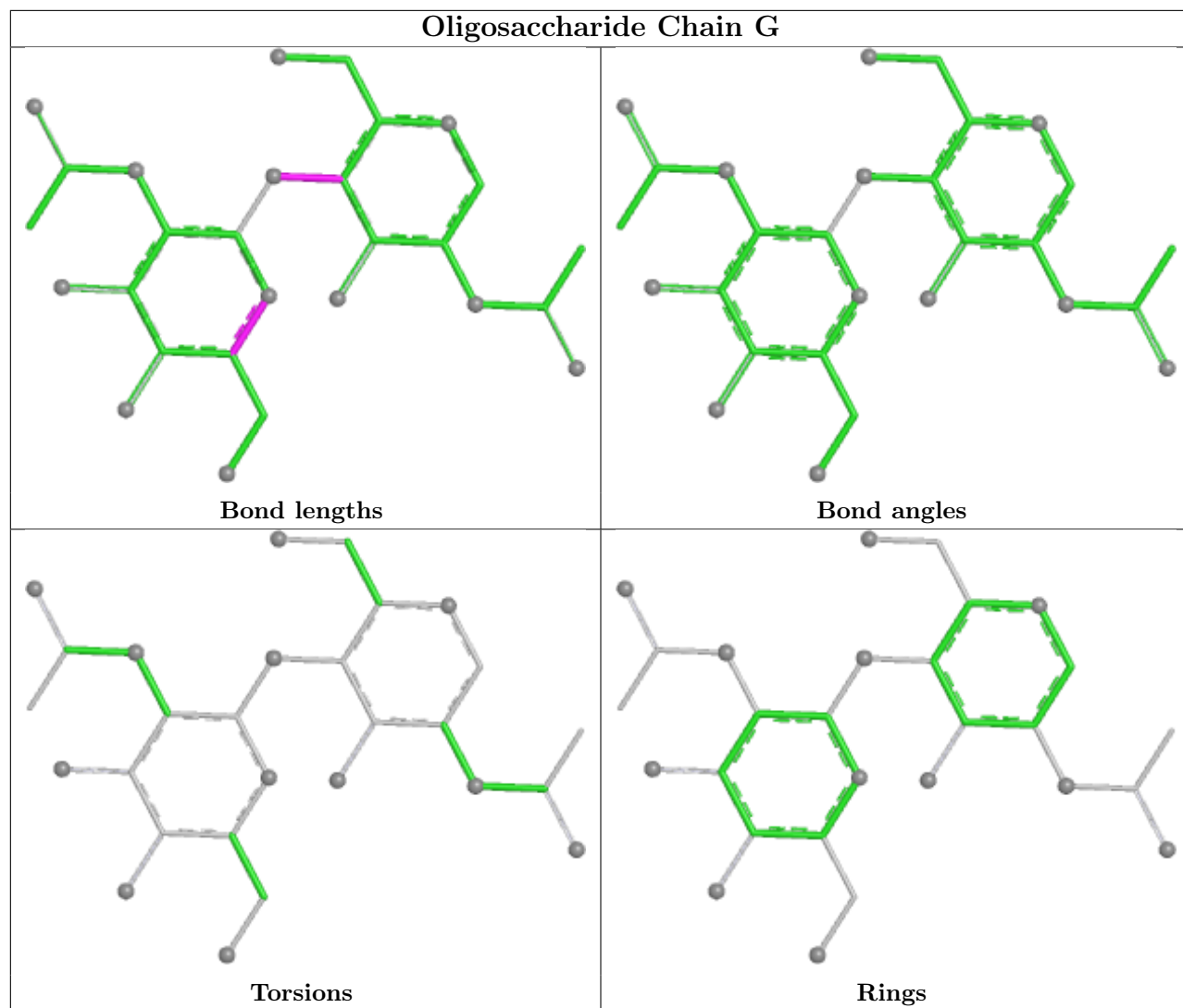
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	c	1	NAG	1	0
4	d	1	NAG	1	0
3	H	1	NAG	4	0
3	H	2	NAG	4	0
3	c	2	NAG	1	0
2	X	1	NAG	2	0
2	M	1	NAG	3	0
4	I	1	NAG	1	0
4	U	1	NAG	1	0

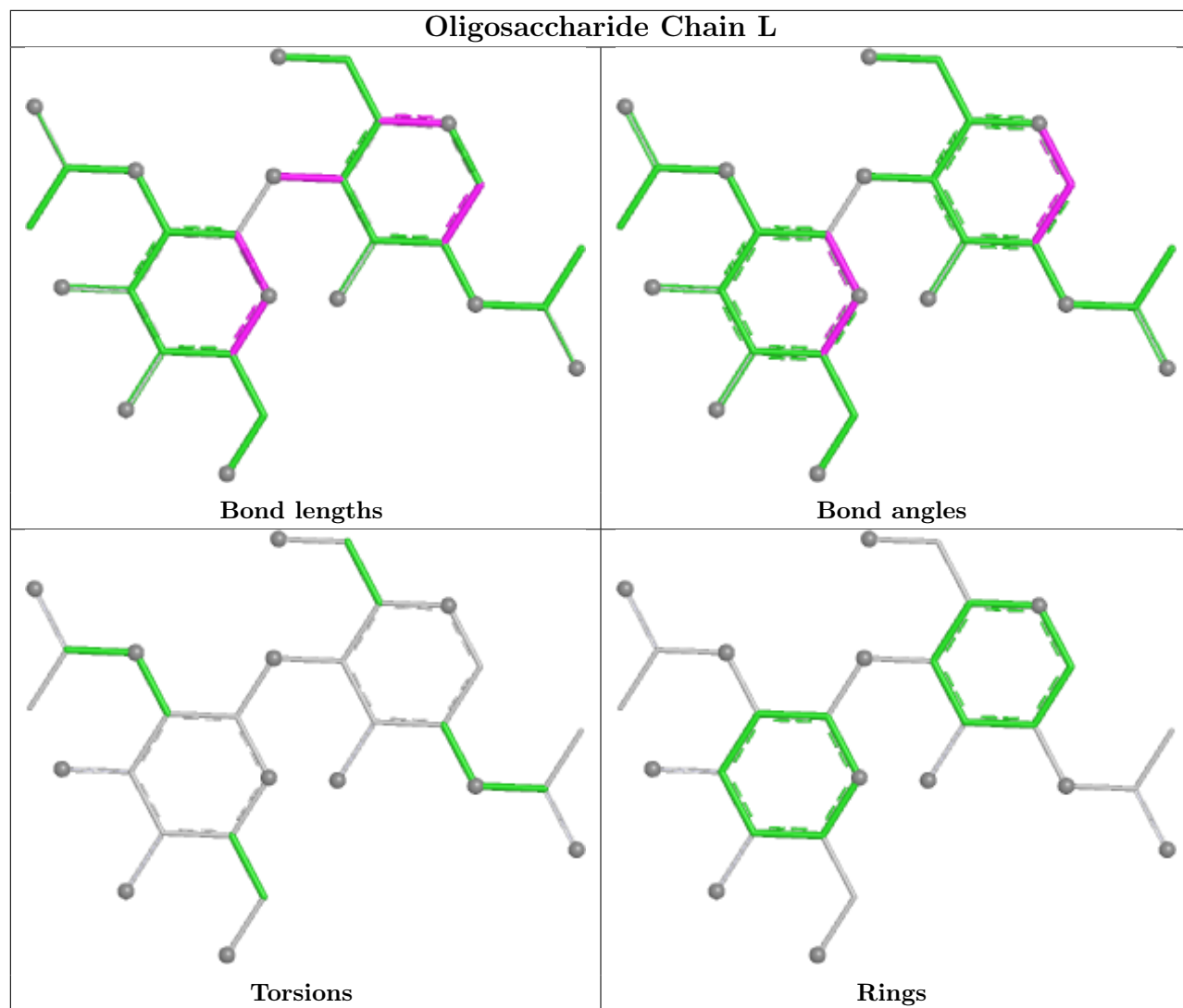
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

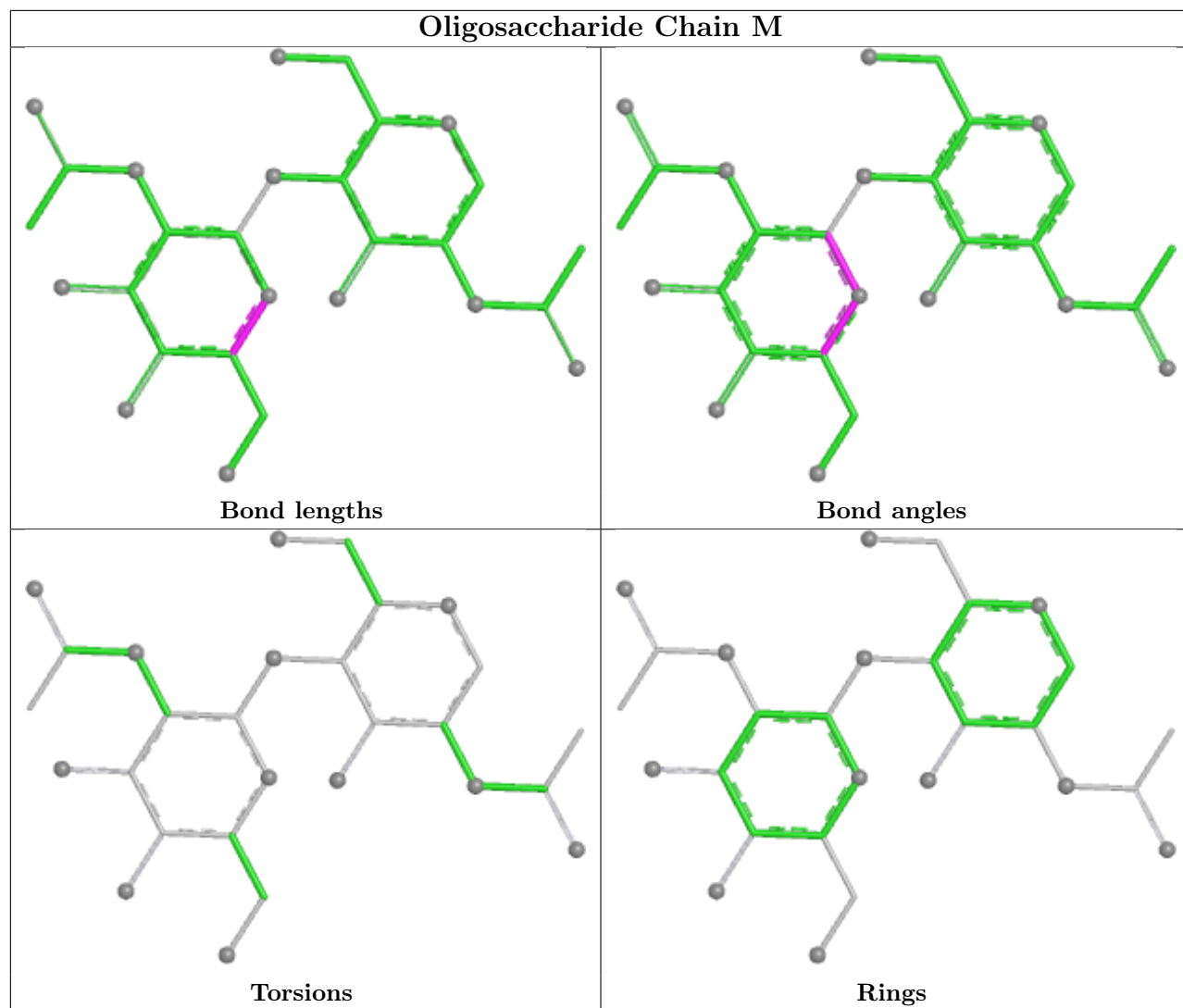


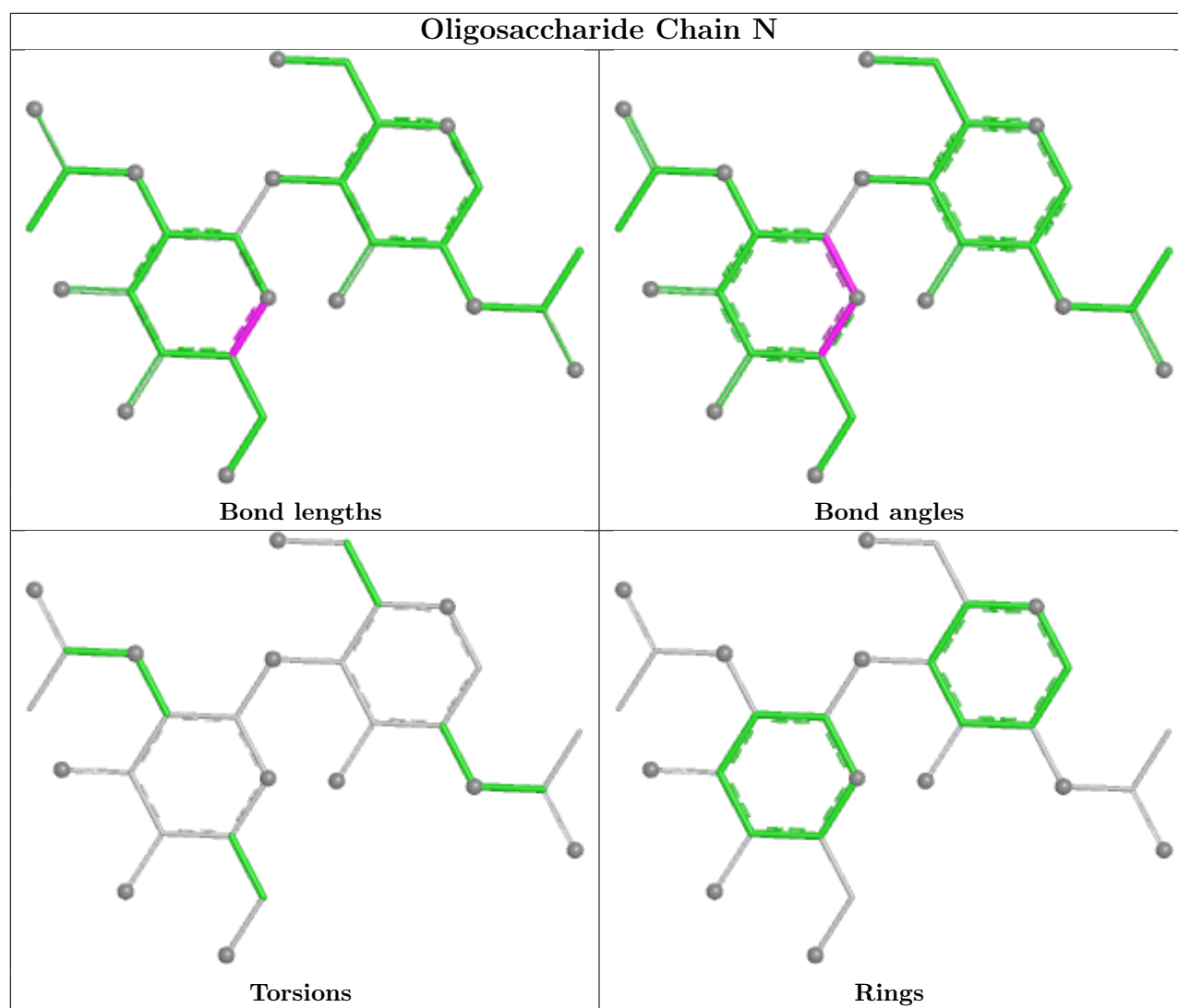


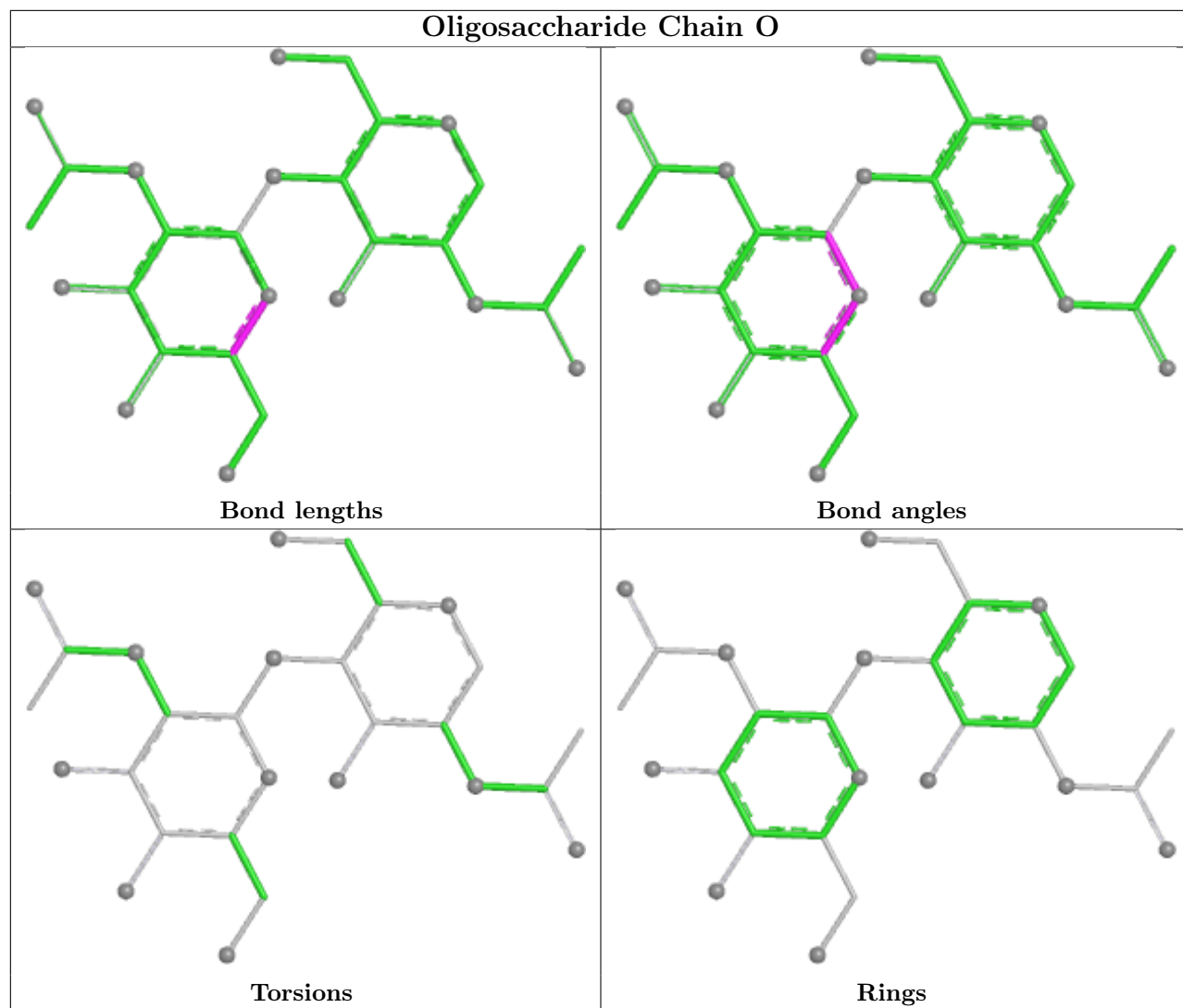


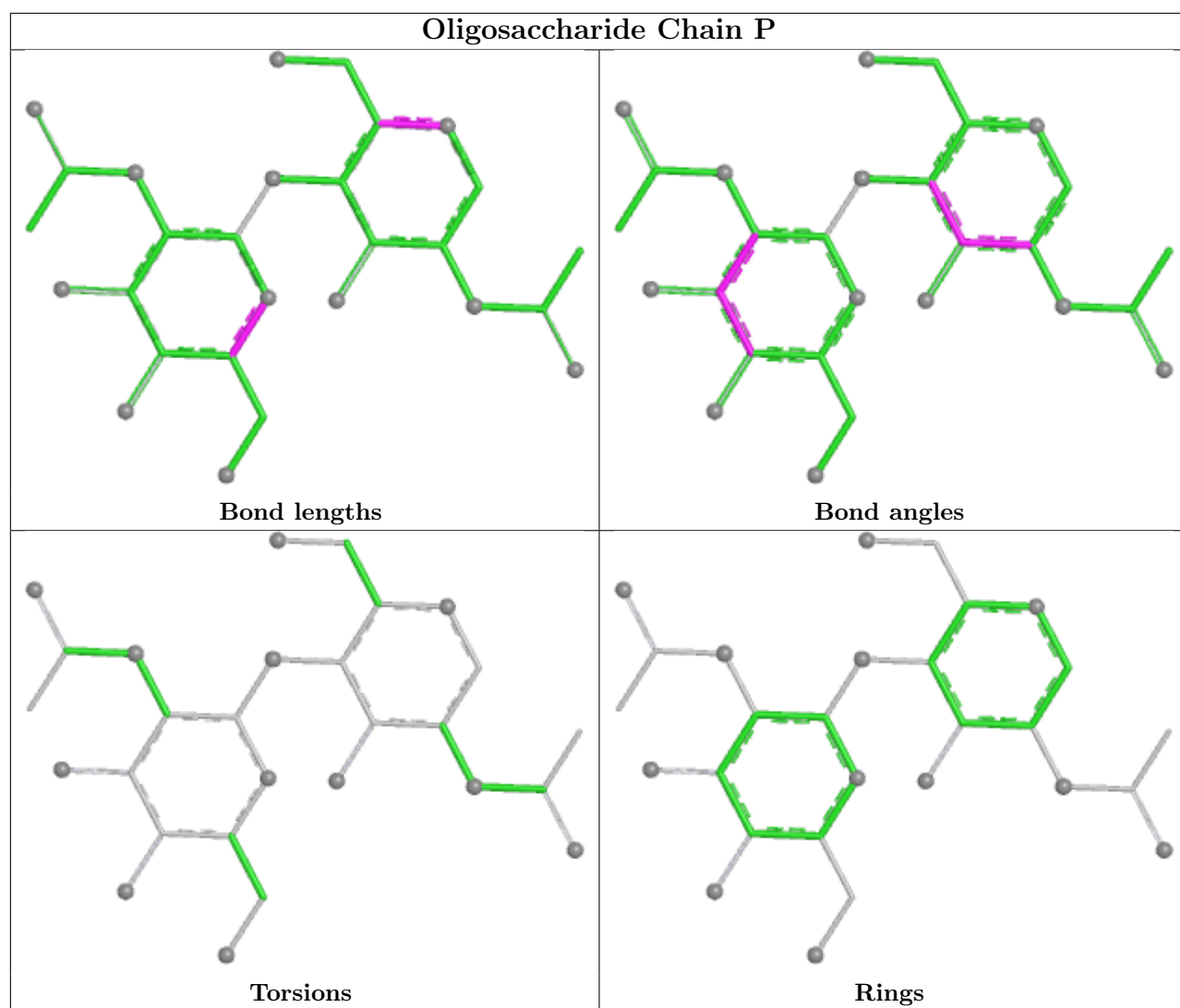


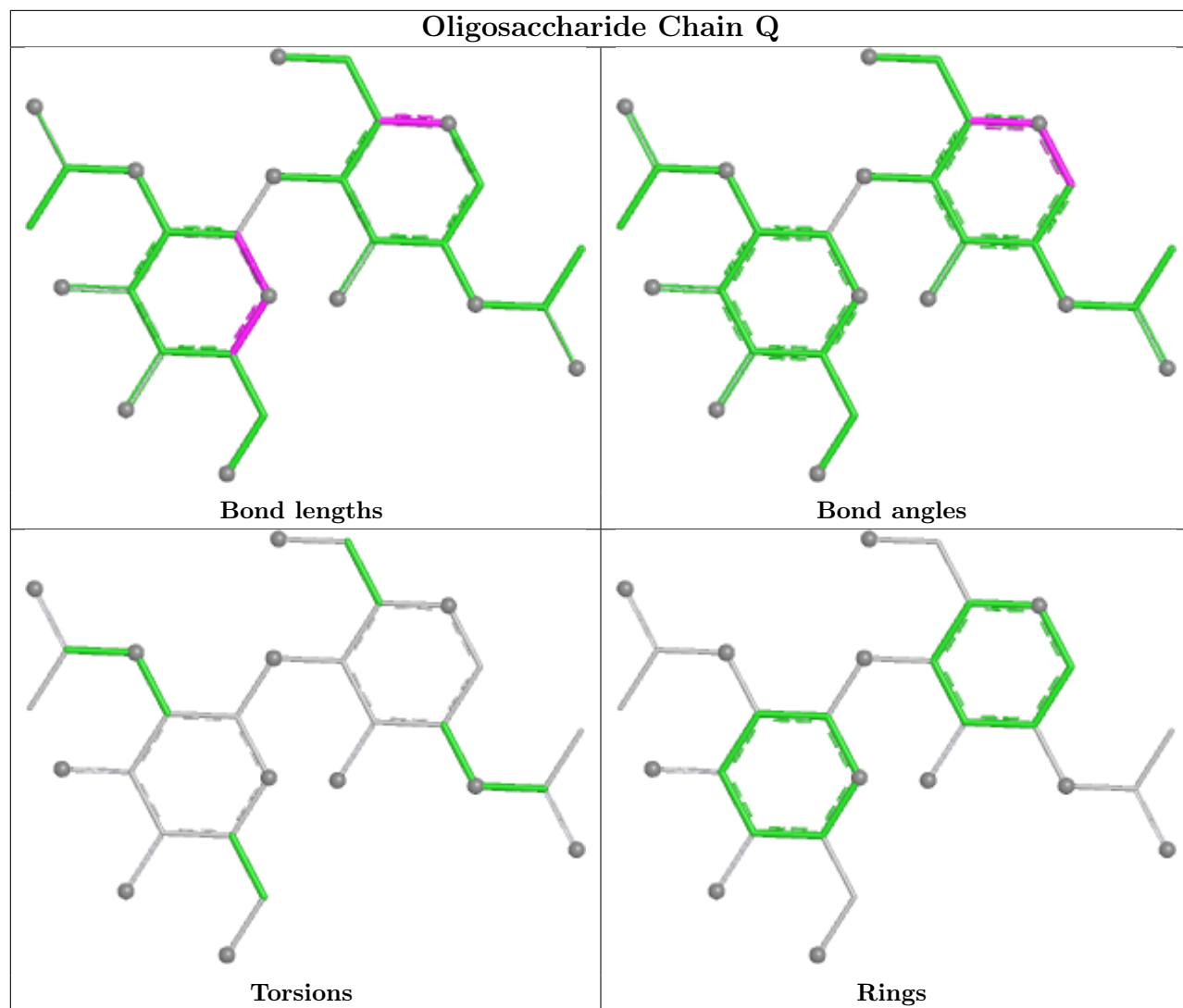


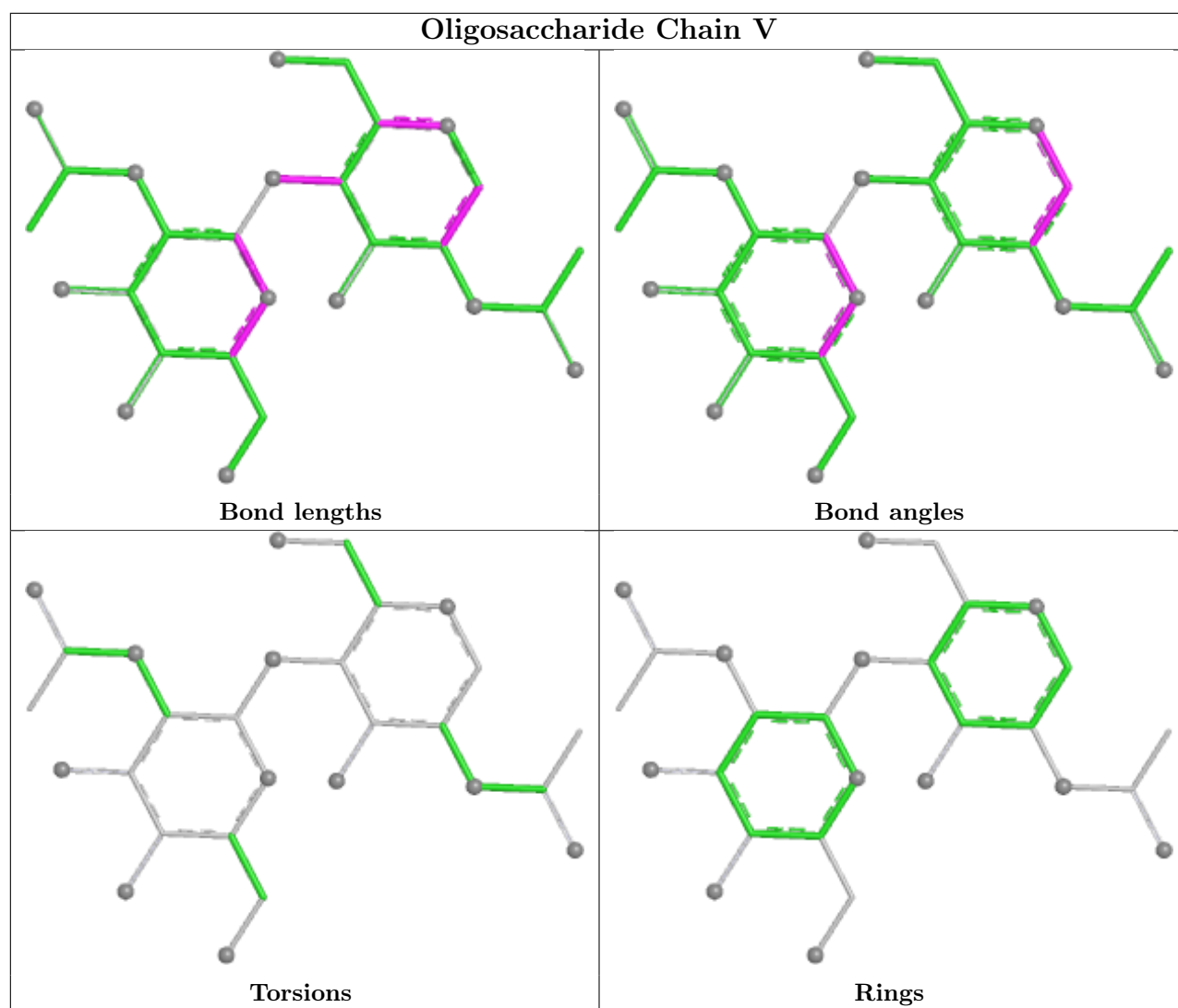


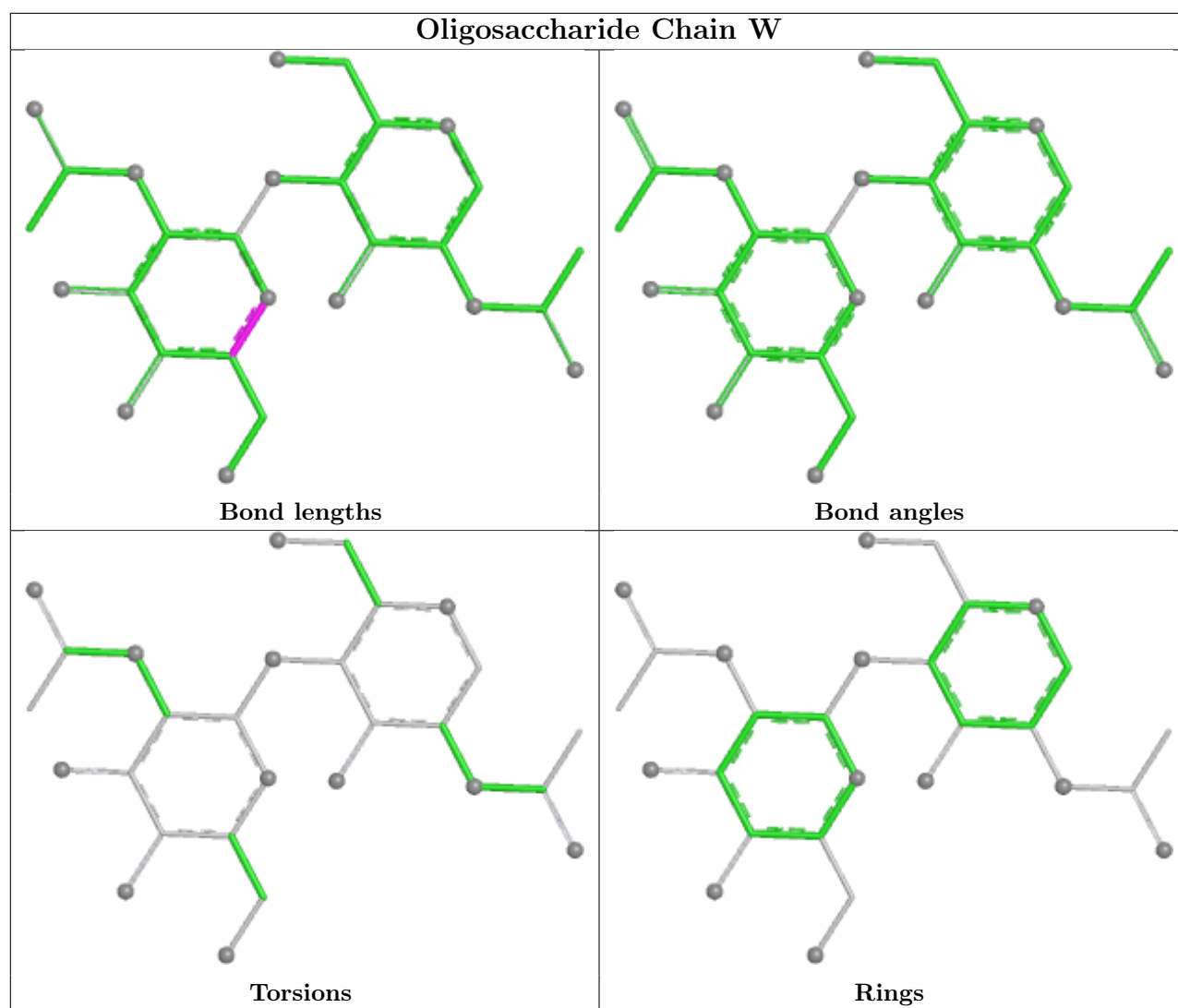


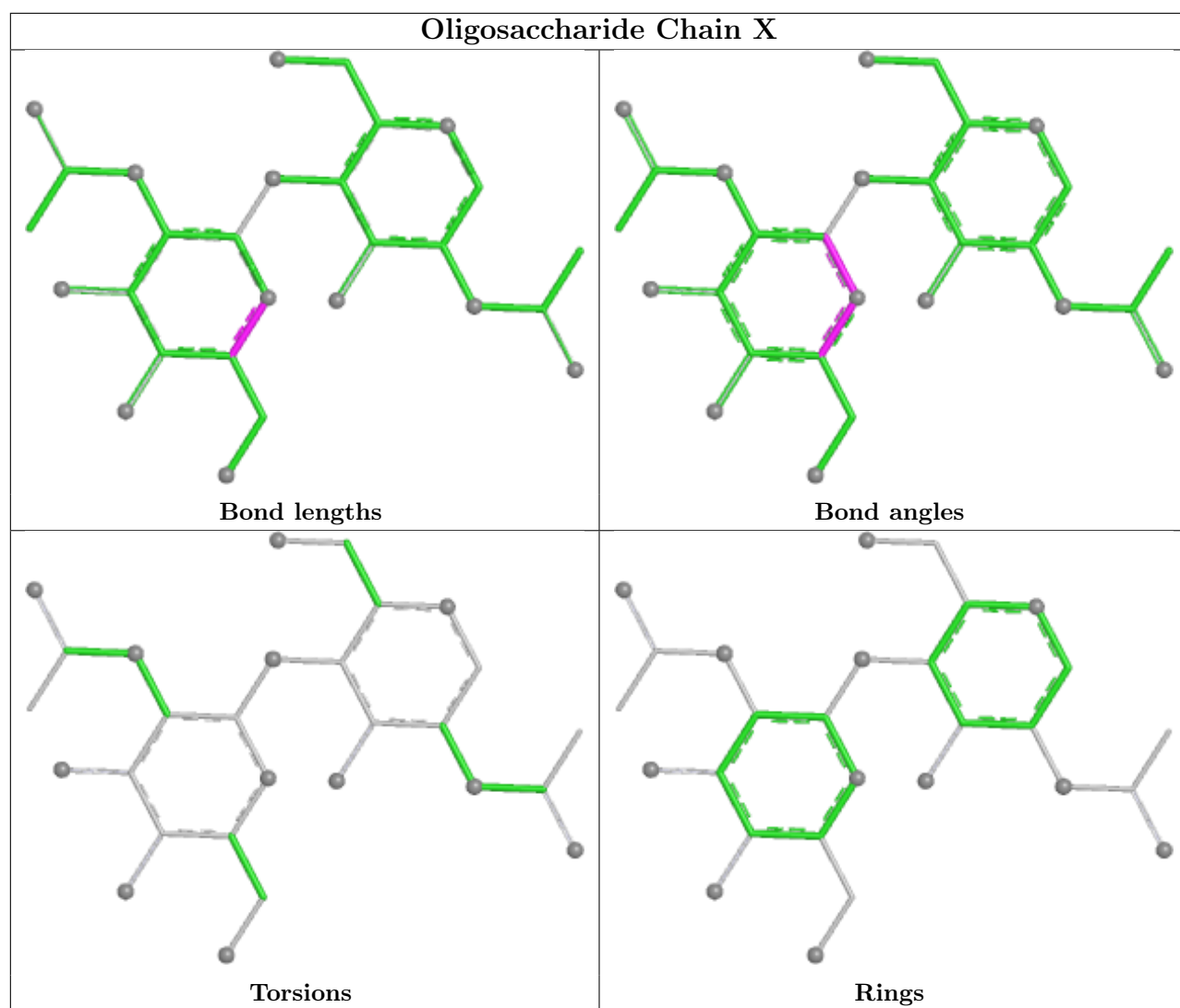


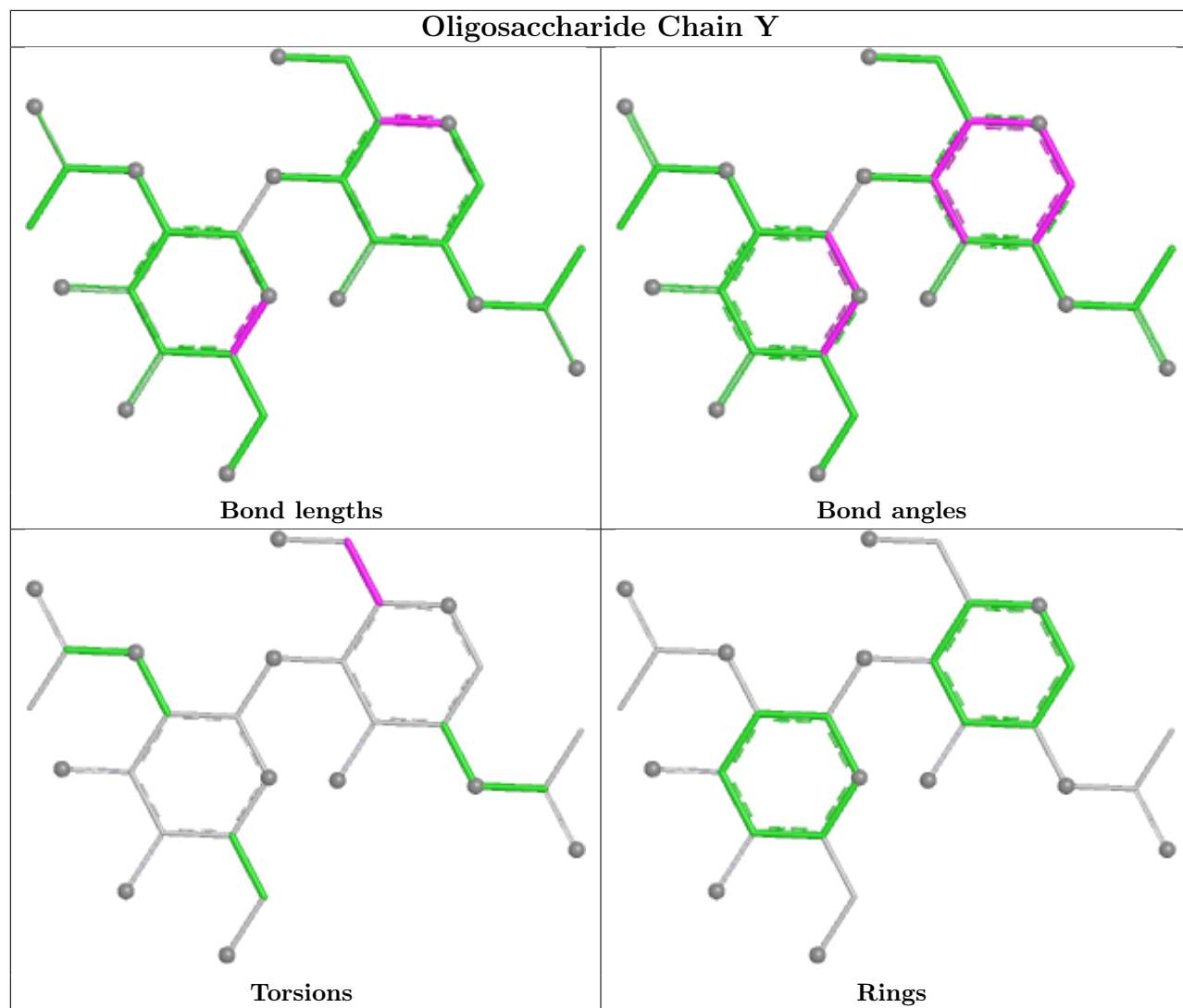


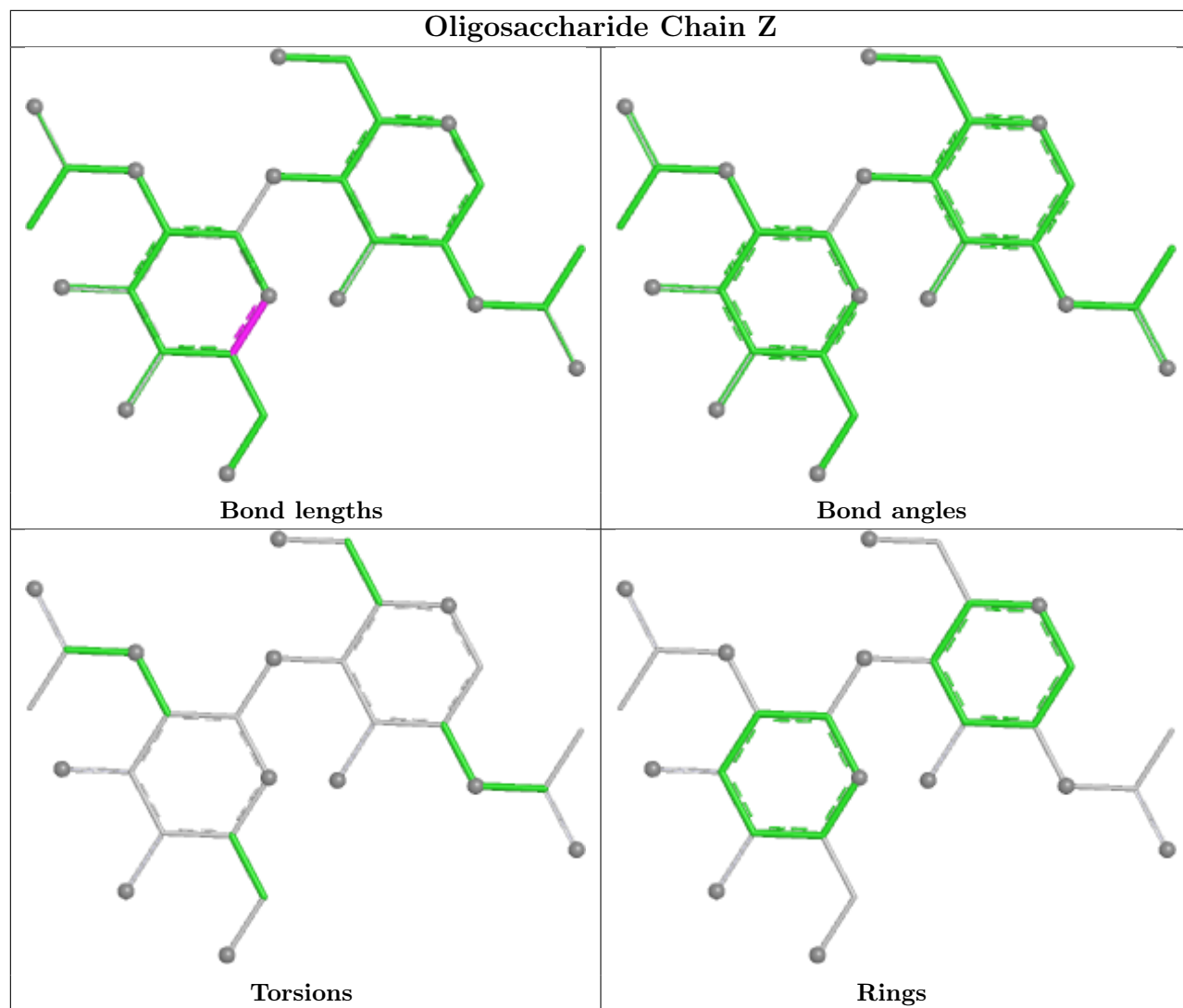


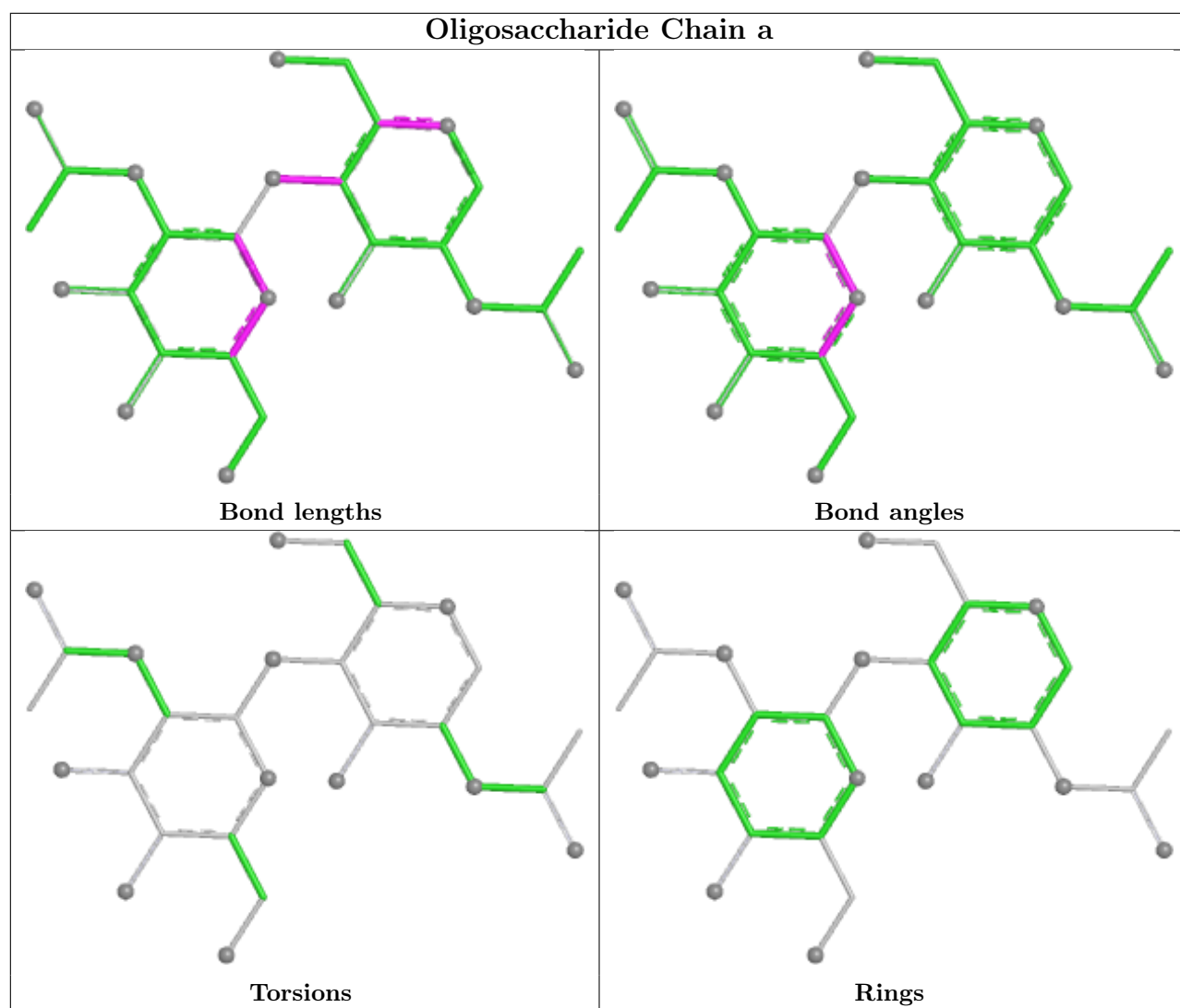


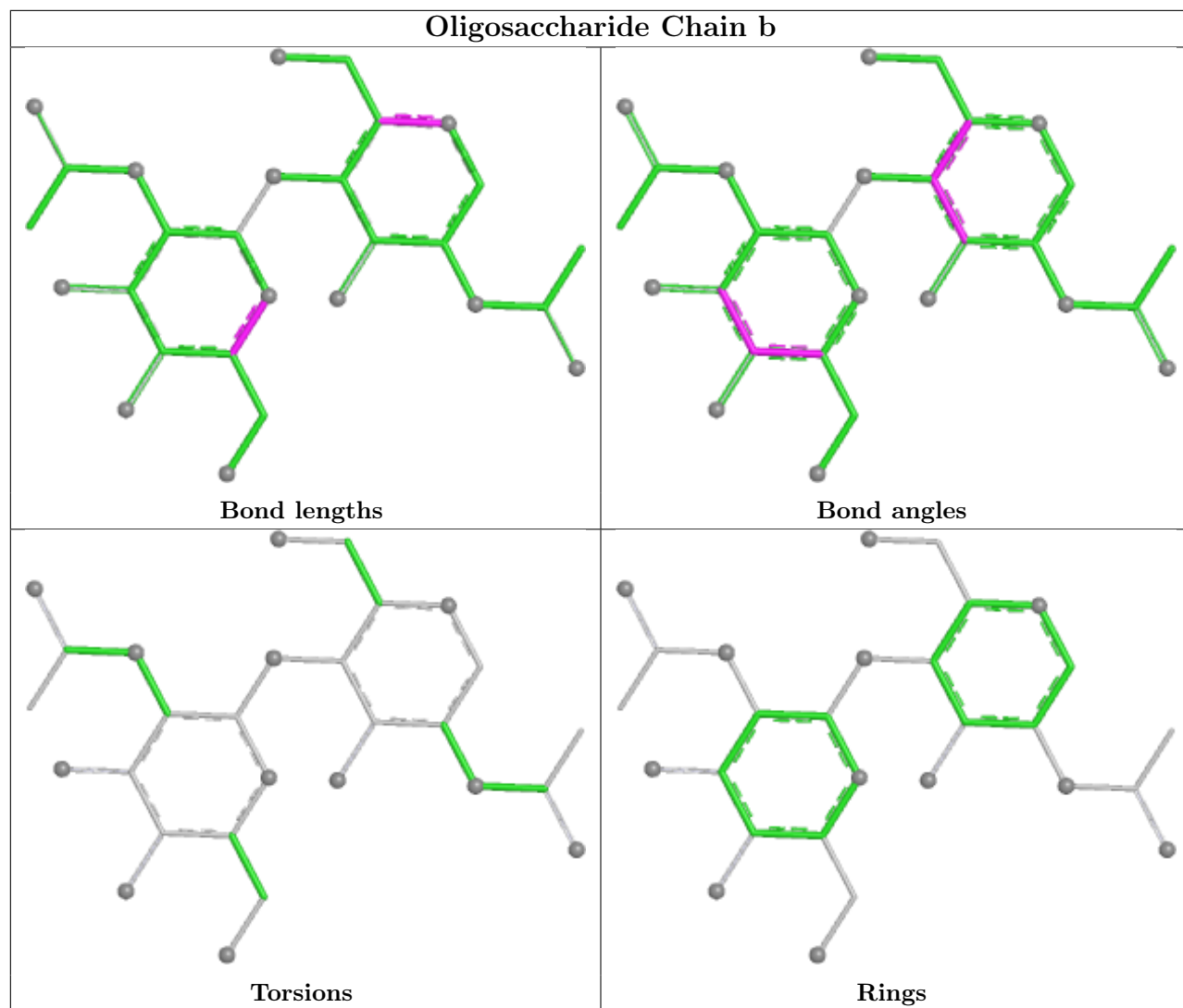


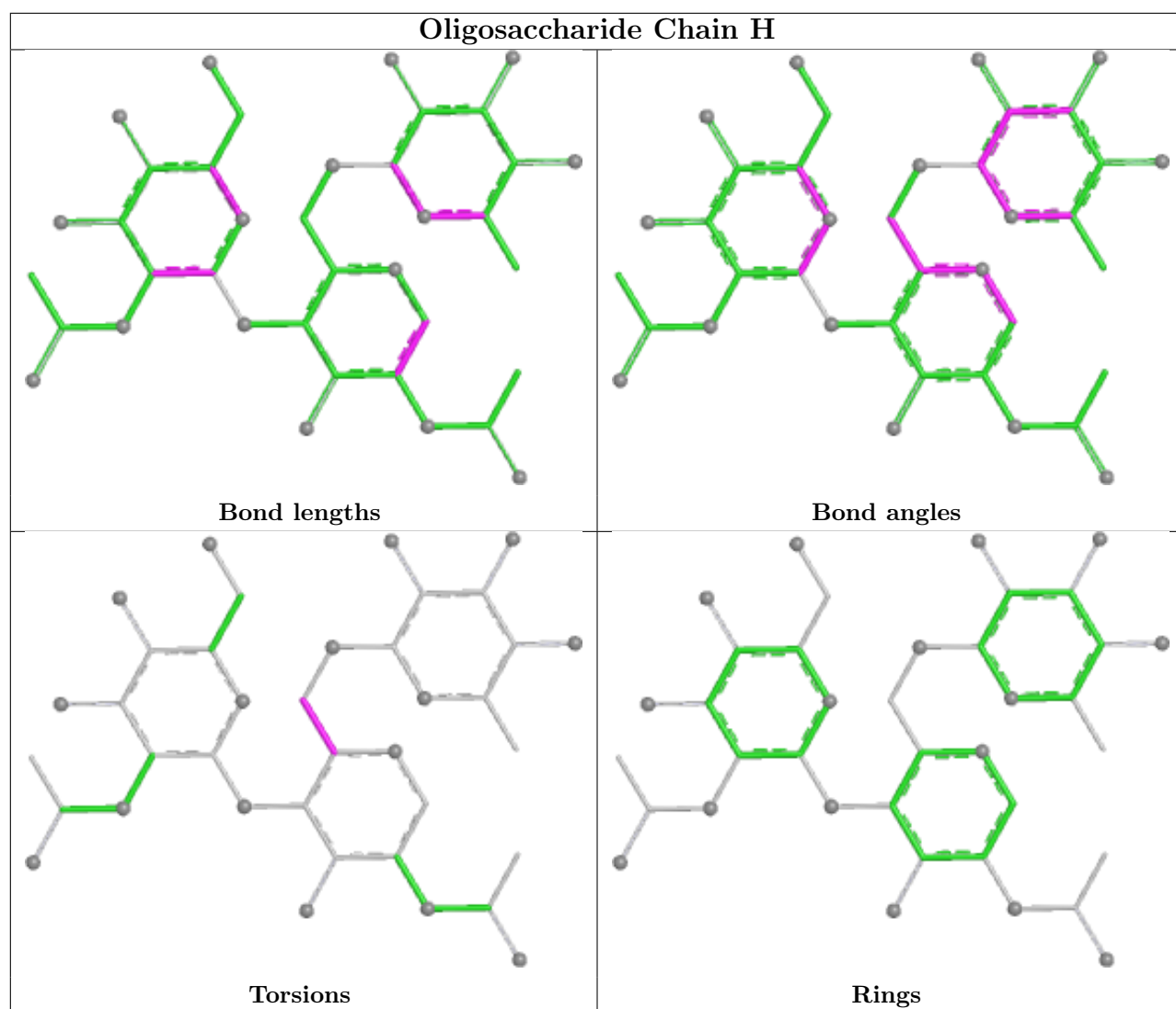


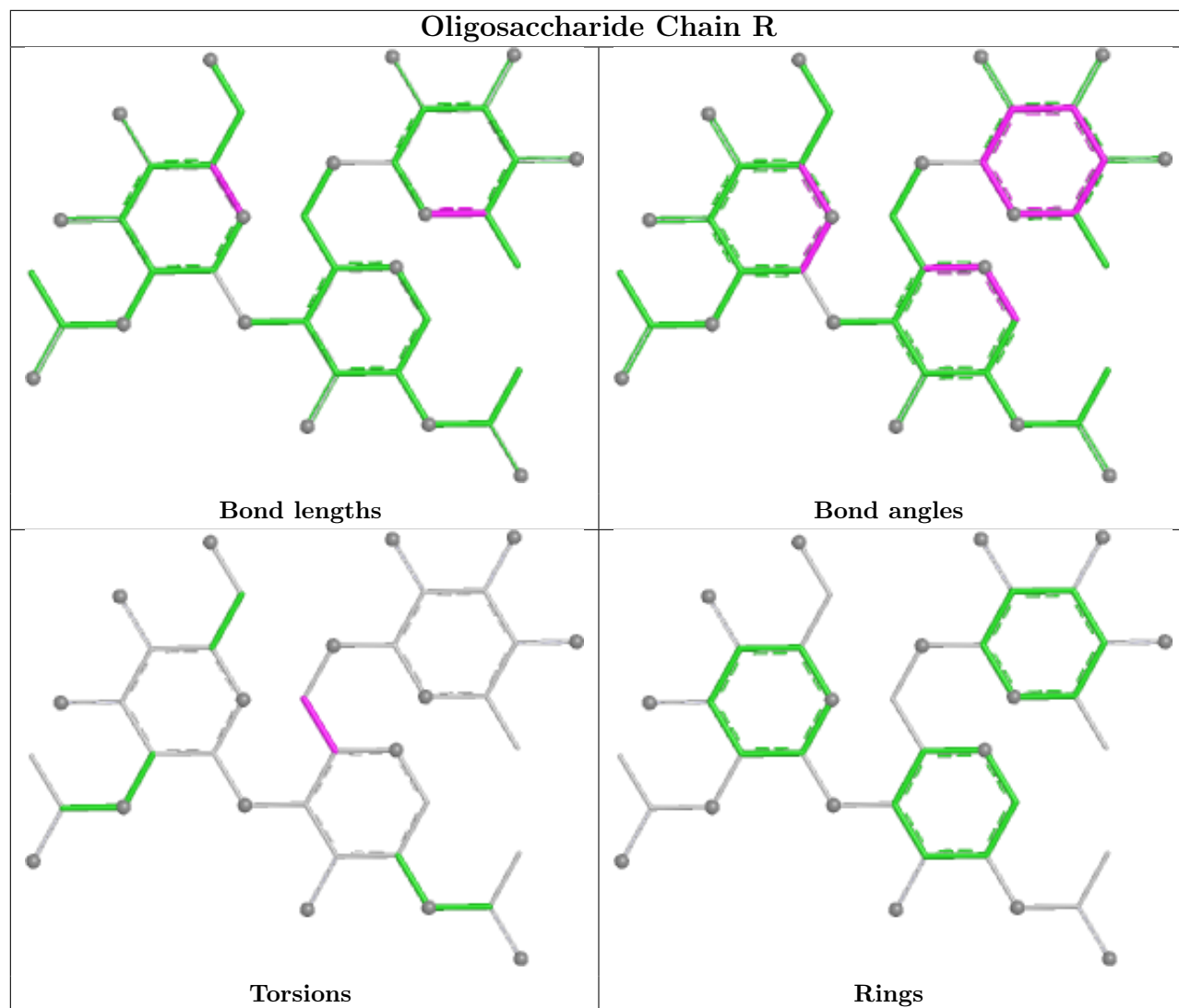


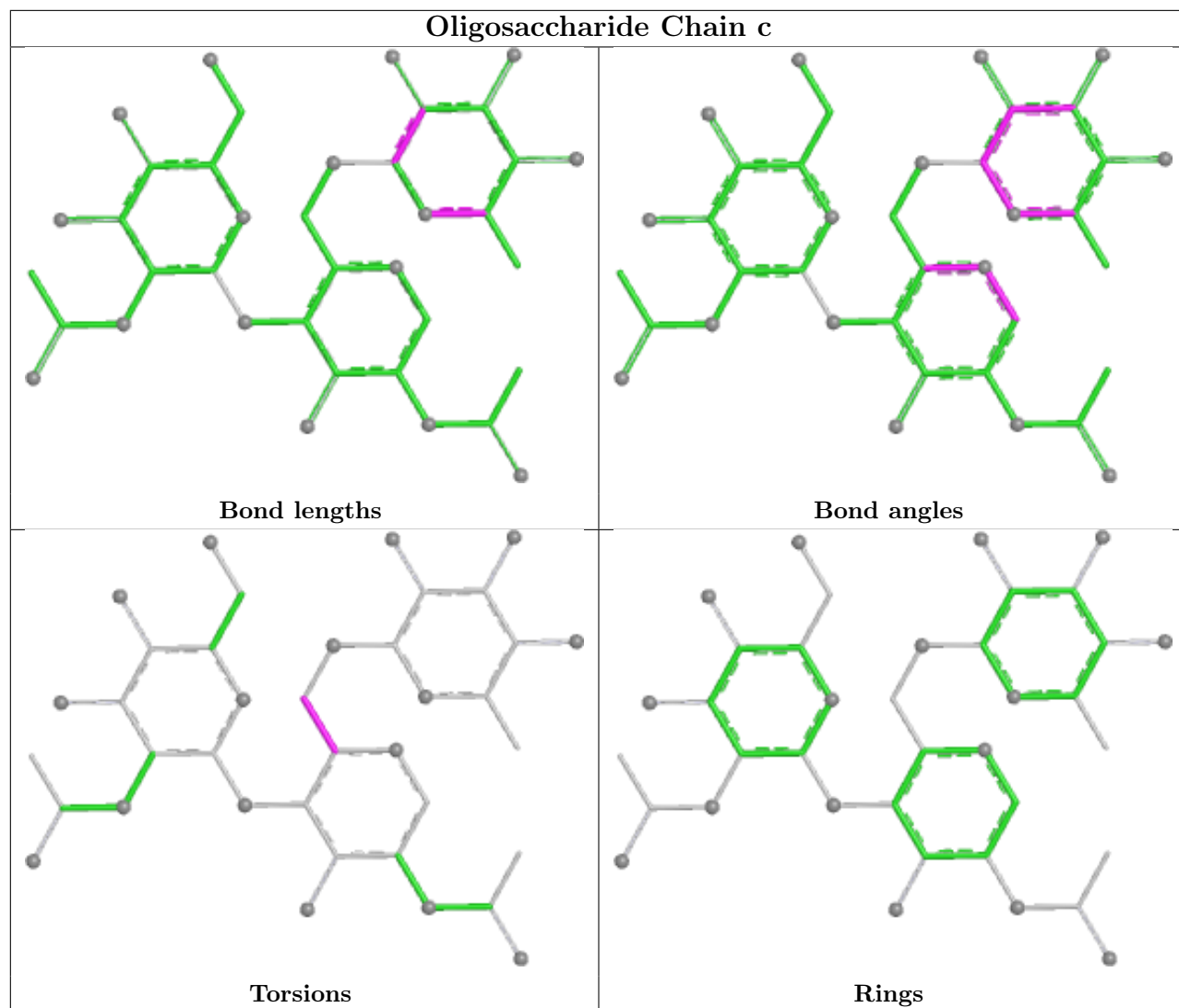


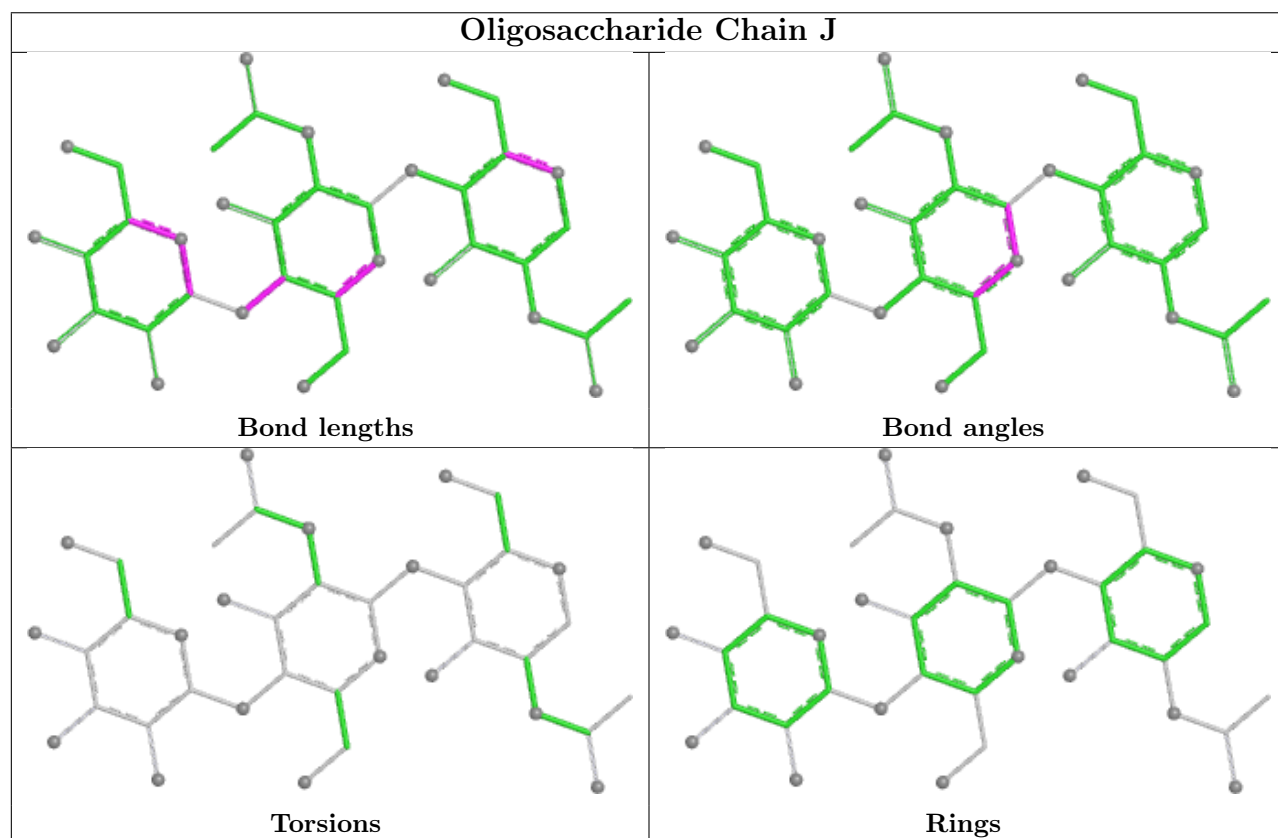
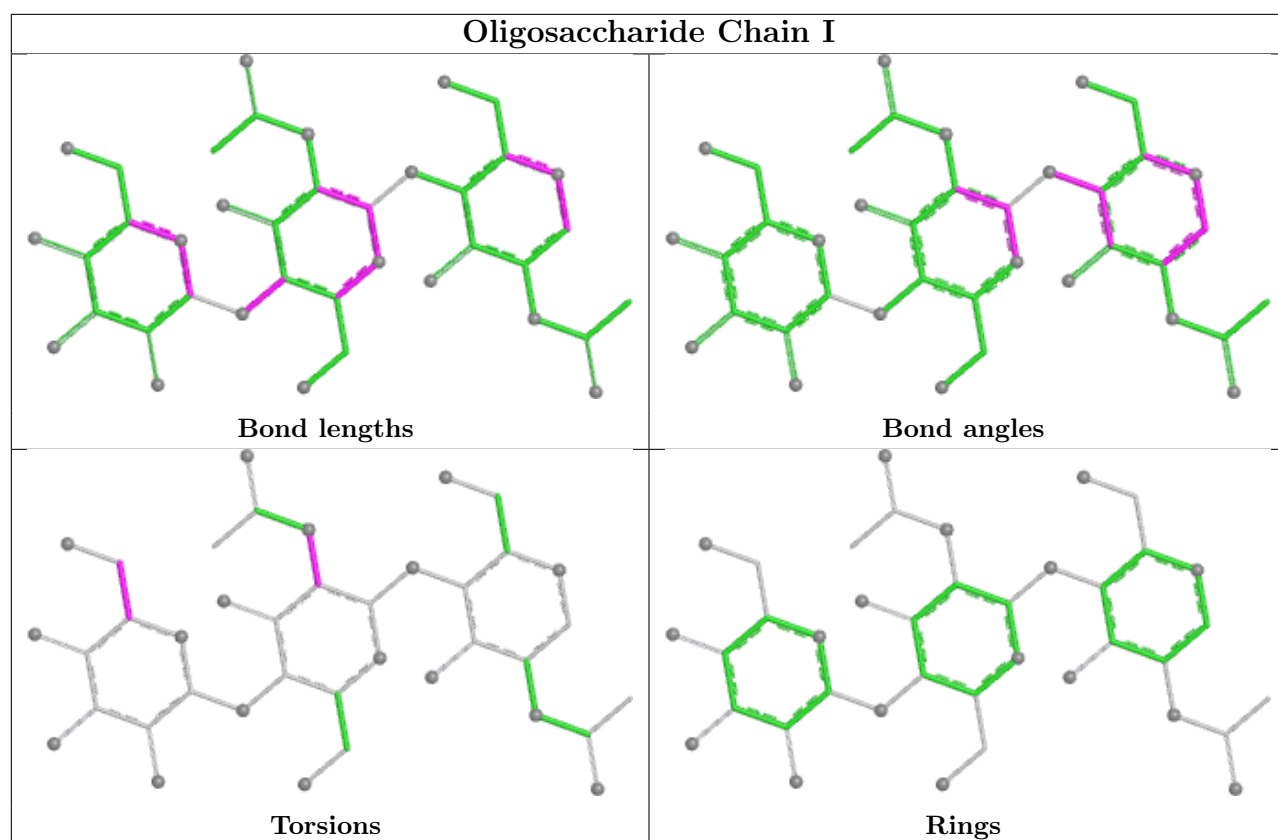


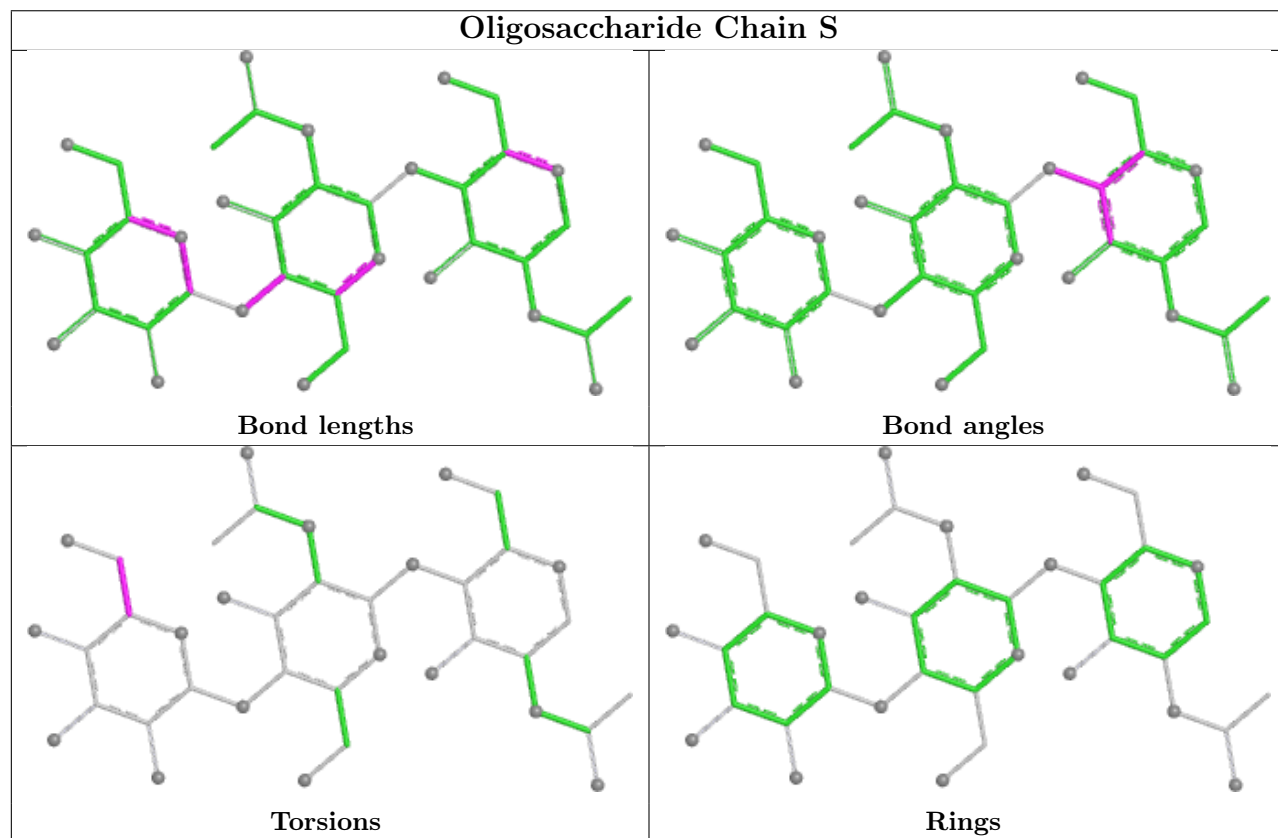
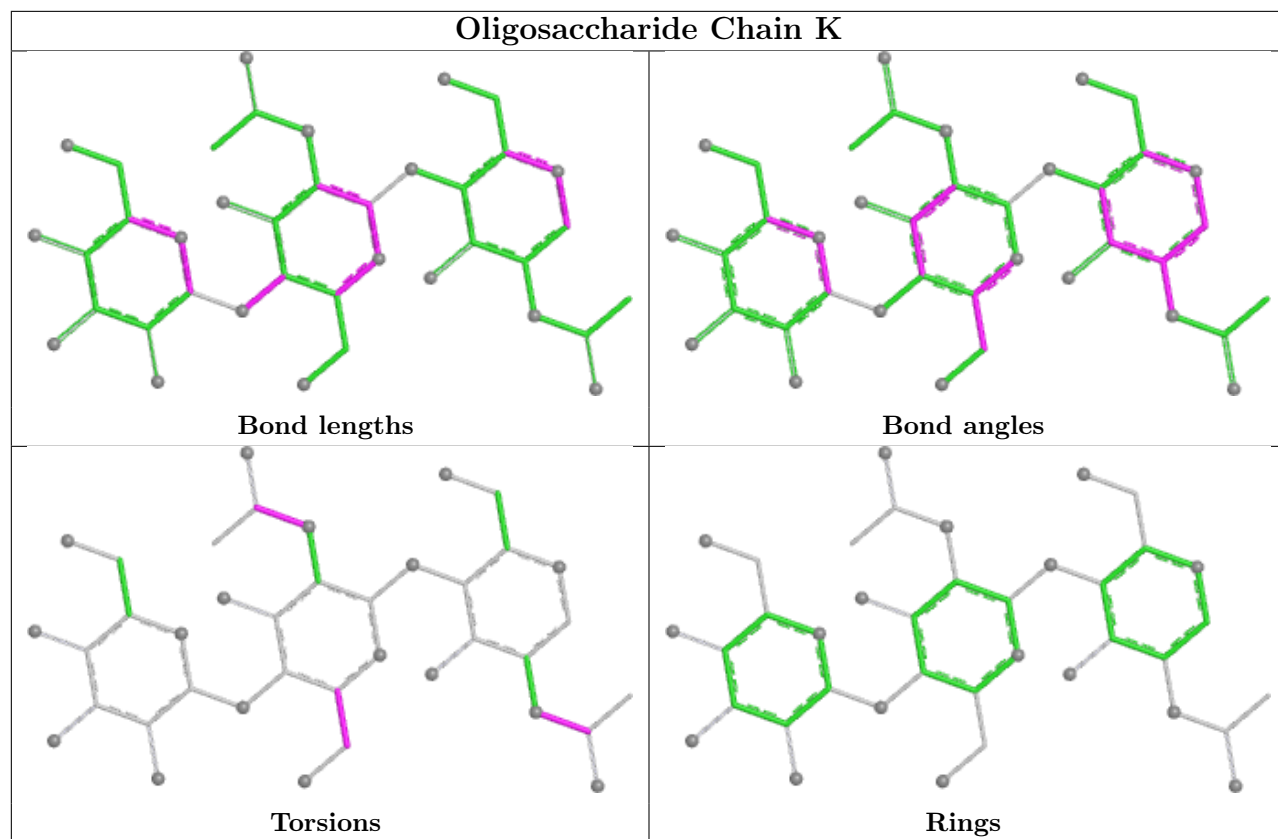


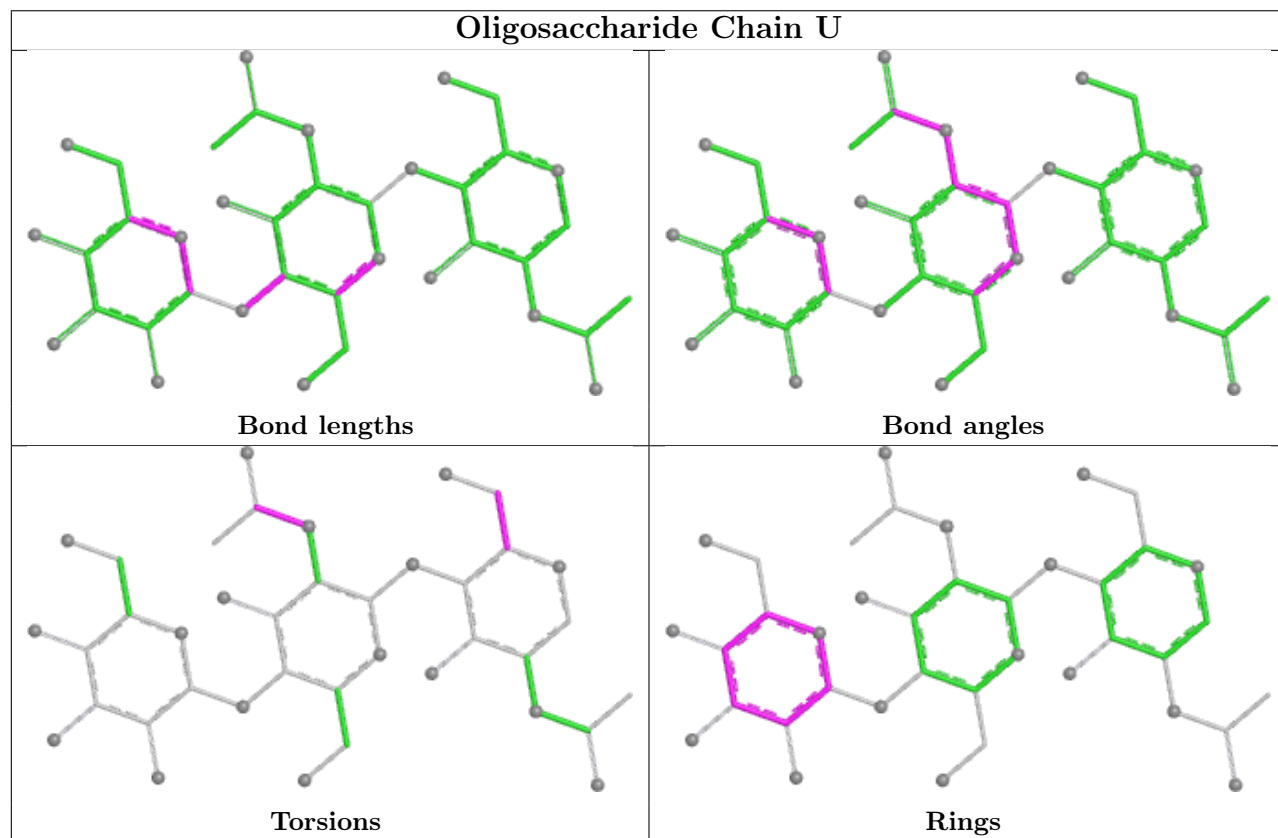
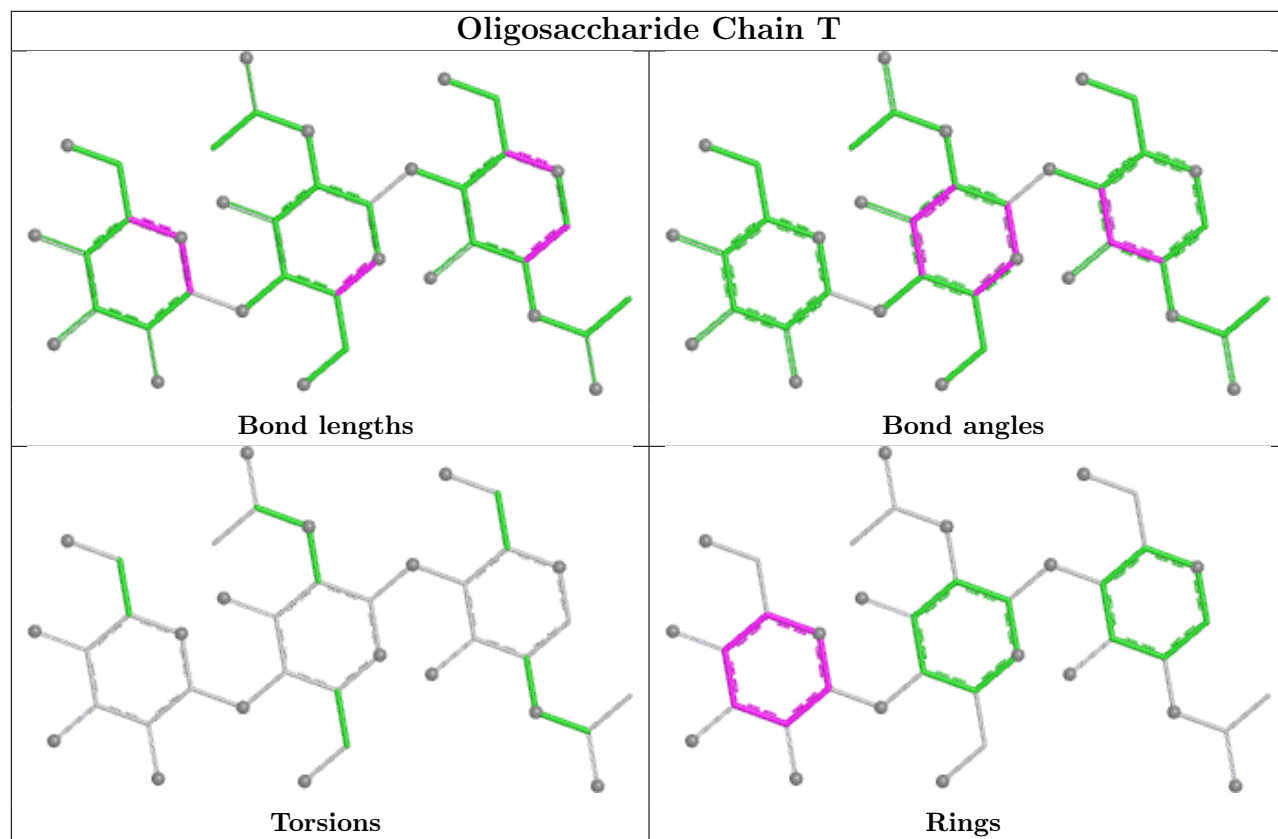


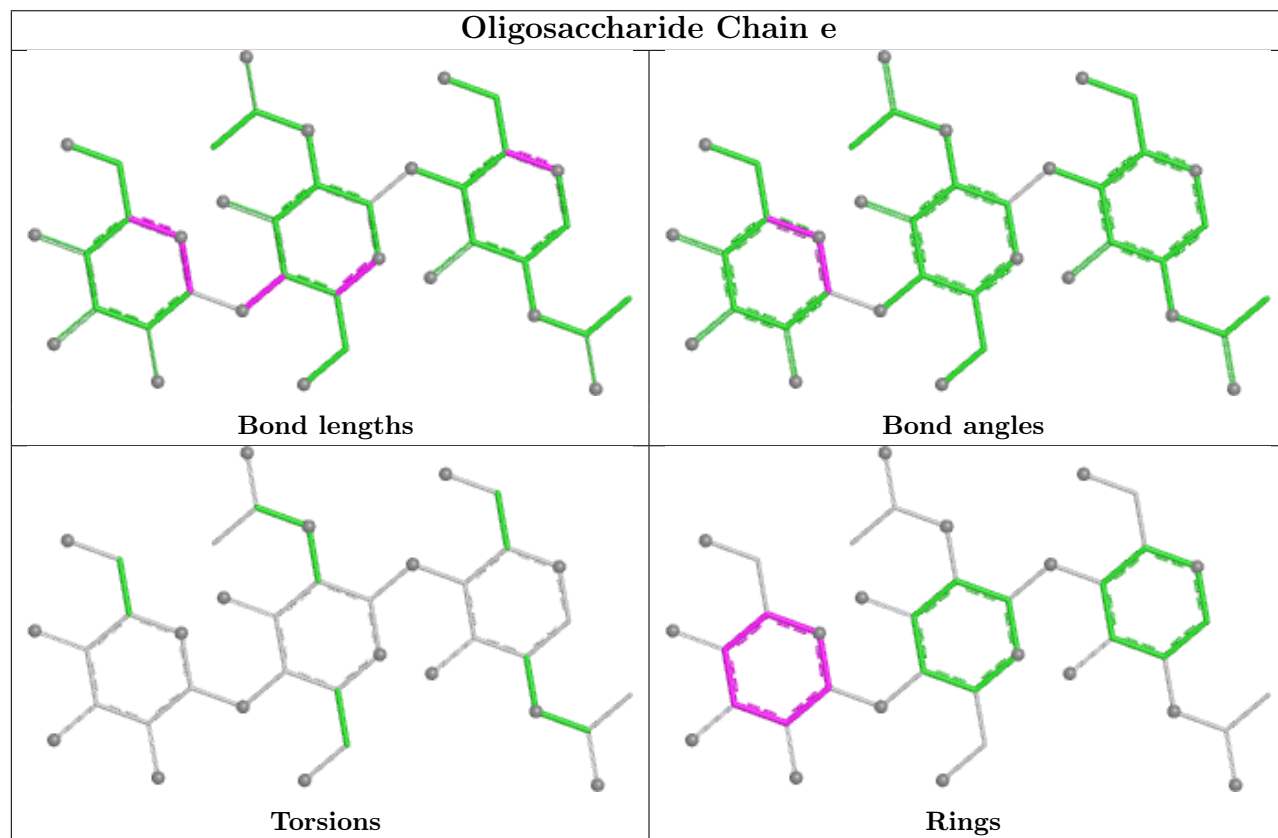
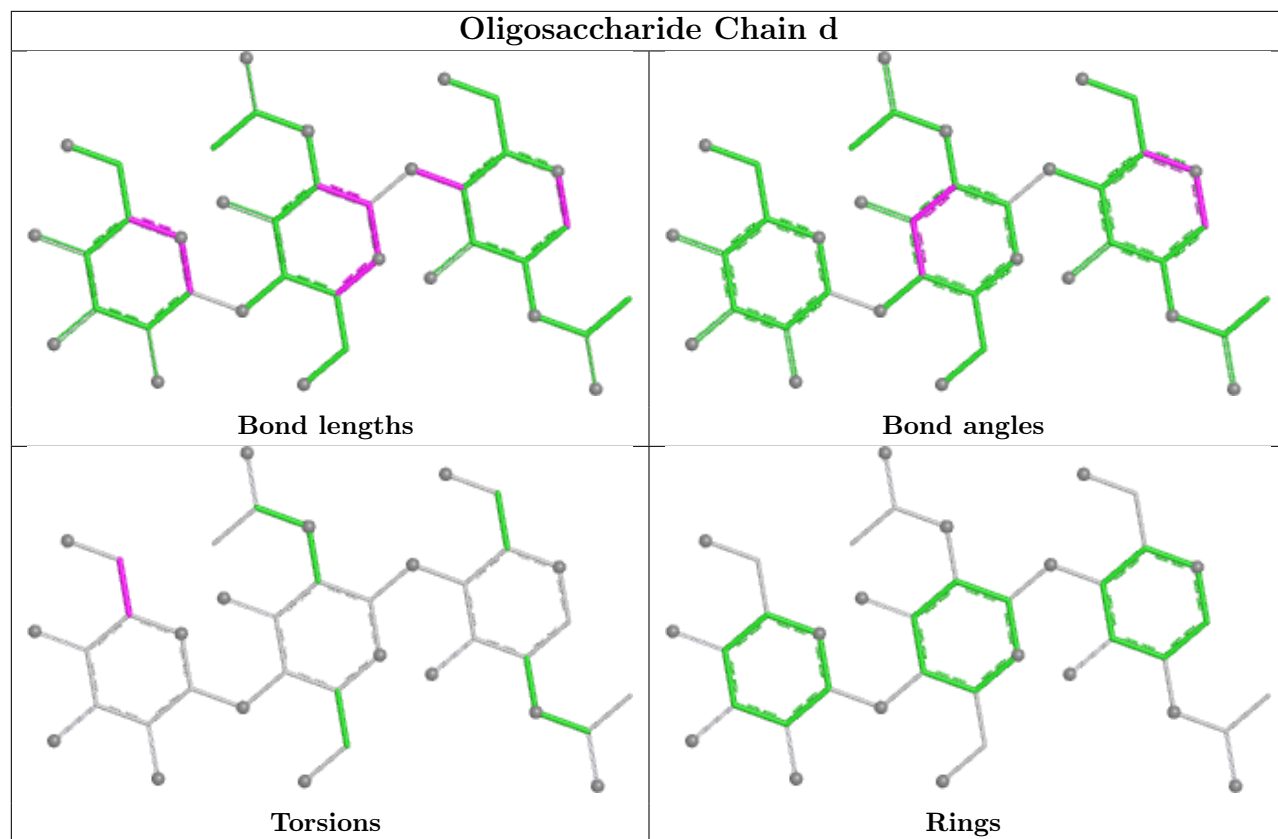


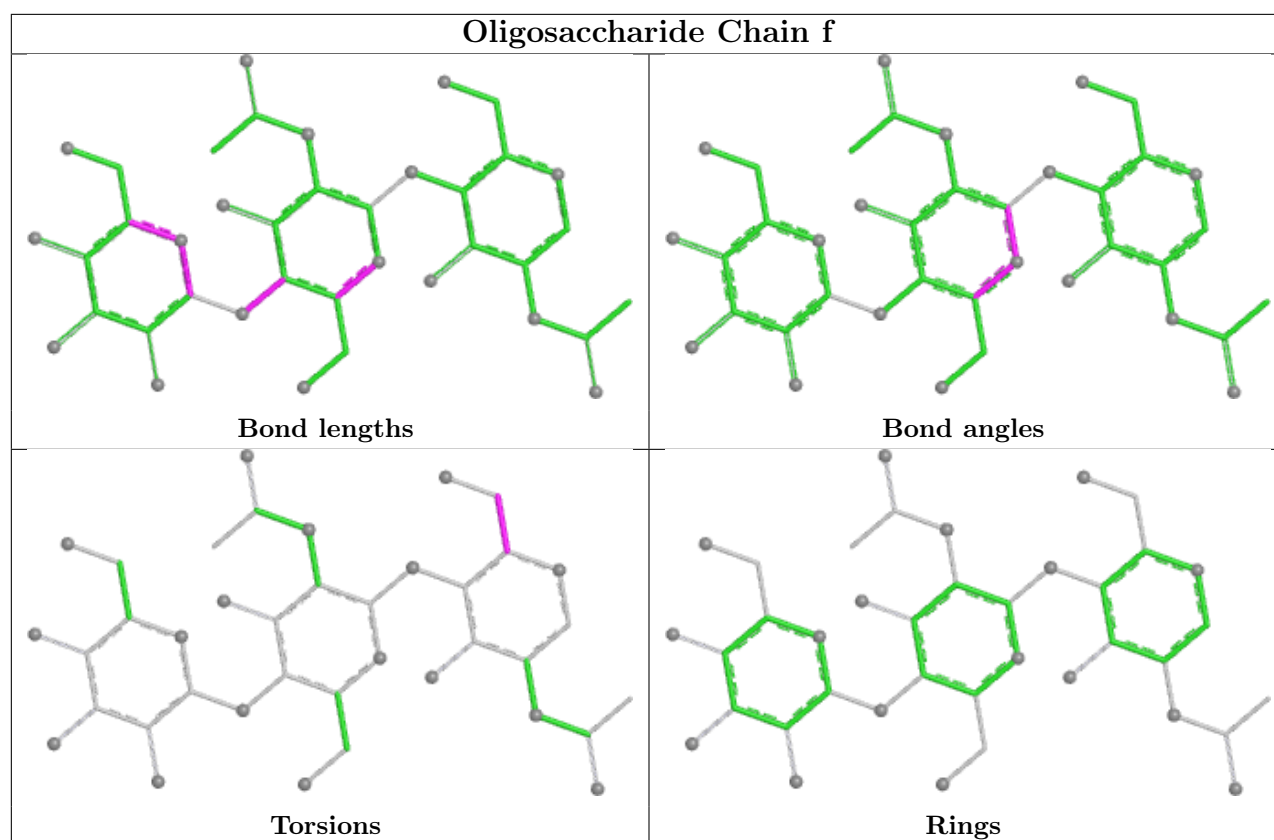












5.6 Ligand geometry [i](#)

28 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	1402	1	14,14,15	1.31	2 (14%)	17,19,21	1.52	4 (23%)
5	NAG	A	1406	1	14,14,15	0.34	0	17,19,21	0.45	0
5	NAG	A	1403	1	14,14,15	0.50	0	17,19,21	0.63	1 (5%)
5	NAG	B	1408	1	14,14,15	1.24	2 (14%)	17,19,21	0.90	1 (5%)
5	NAG	B	1401	1	14,14,15	1.36	1 (7%)	17,19,21	1.46	3 (17%)
5	NAG	B	1402	1	14,14,15	0.50	0	17,19,21	0.70	1 (5%)
5	NAG	B	1403	1	14,14,15	1.14	1 (7%)	17,19,21	0.96	1 (5%)
5	NAG	B	1405	1	14,14,15	1.18	2 (14%)	17,19,21	0.83	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	A	1407	1	14,14,15	0.32	0	17,19,21	0.45	0
5	NAG	B	1406	1	14,14,15	1.16	1 (7%)	17,19,21	1.16	2 (11%)
5	NAG	A	1409	1	14,14,15	0.29	0	17,19,21	0.40	0
5	NAG	C	1401	1	14,14,15	0.61	0	17,19,21	0.44	0
5	NAG	C	1403	1	14,14,15	1.11	1 (7%)	17,19,21	1.10	1 (5%)
5	NAG	C	1405	1	14,14,15	0.23	0	17,19,21	0.44	0
5	NAG	B	1409	1	14,14,15	1.27	1 (7%)	17,19,21	0.63	0
5	NAG	C	1408	1	14,14,15	1.29	2 (14%)	17,19,21	0.80	1 (5%)
5	NAG	C	1402	1	14,14,15	1.31	2 (14%)	17,19,21	0.61	0
5	NAG	B	1407	1	14,14,15	0.22	0	17,19,21	0.45	0
5	NAG	C	1407	1	14,14,15	1.51	3 (21%)	17,19,21	0.65	0
5	NAG	B	1404	1	14,14,15	1.14	1 (7%)	17,19,21	1.11	1 (5%)
5	NAG	A	1411	1	14,14,15	0.29	0	17,19,21	0.42	0
5	NAG	C	1404	1	14,14,15	1.65	2 (14%)	17,19,21	0.68	0
5	NAG	A	1408	1	14,14,15	1.08	1 (7%)	17,19,21	0.70	0
5	NAG	A	1410	1	14,14,15	1.12	1 (7%)	17,19,21	1.18	2 (11%)
5	NAG	A	1401	1	14,14,15	1.34	2 (14%)	17,19,21	0.95	1 (5%)
5	NAG	A	1405	1	14,14,15	1.17	1 (7%)	17,19,21	0.76	0
5	NAG	C	1406	1	14,14,15	1.04	1 (7%)	17,19,21	0.75	0
5	NAG	A	1404	1	14,14,15	1.03	1 (7%)	17,19,21	1.00	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1402	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1406	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1403	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1408	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1401	1	-	1/6/23/26	0/1/1/1
5	NAG	B	1402	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1403	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1405	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	1409	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	C	1401	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1403	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1405	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1409	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1408	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	1407	1	-	1/6/23/26	0/1/1/1
5	NAG	C	1407	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1404	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1411	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	1408	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1410	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	-	0/6/23/26	0/1/1/1
5	NAG	C	1406	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1404	1	-	0/6/23/26	0/1/1/1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	1404	NAG	C1-C2	4.65	1.58	1.52
5	B	1404	NAG	O5-C1	-4.17	1.36	1.43
5	C	1403	NAG	O5-C1	-4.08	1.36	1.43
5	B	1401	NAG	O5-C5	3.86	1.51	1.43
5	A	1405	NAG	O5-C5	3.29	1.49	1.43
5	A	1402	NAG	O5-C5	2.99	1.49	1.43
5	C	1407	NAG	C8-C7	2.99	1.56	1.50
5	A	1401	NAG	C1-C2	2.97	1.56	1.52
5	C	1408	NAG	O5-C5	2.92	1.49	1.43
5	C	1407	NAG	O5-C5	2.91	1.49	1.43
5	C	1406	NAG	O5-C5	2.77	1.48	1.43
5	A	1404	NAG	O5-C5	2.70	1.48	1.43
5	B	1405	NAG	O5-C5	2.64	1.48	1.43
5	C	1402	NAG	O5-C5	2.56	1.48	1.43
5	B	1409	NAG	O5-C5	2.55	1.48	1.43
5	A	1410	NAG	O5-C5	2.50	1.48	1.43
5	B	1403	NAG	O5-C5	2.45	1.48	1.43
5	B	1408	NAG	O5-C5	2.40	1.48	1.43
5	C	1404	NAG	O5-C5	2.36	1.48	1.43
5	B	1408	NAG	C1-C2	2.29	1.55	1.52
5	C	1402	NAG	C1-C2	2.28	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1401	NAG	O5-C5	2.24	1.47	1.43
5	C	1407	NAG	C1-C2	2.15	1.55	1.52
5	B	1406	NAG	O5-C5	2.12	1.47	1.43
5	B	1405	NAG	C1-C2	2.11	1.55	1.52
5	C	1408	NAG	C1-C2	2.11	1.55	1.52
5	A	1408	NAG	O5-C5	2.06	1.47	1.43
5	A	1402	NAG	C1-C2	2.02	1.55	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1401	NAG	C1-O5-C5	4.20	117.82	112.19
5	A	1410	NAG	C1-O5-C5	3.48	116.85	112.19
5	A	1402	NAG	C1-C2-N2	3.18	115.44	110.43
5	B	1408	NAG	C1-O5-C5	3.06	116.29	112.19
5	A	1404	NAG	C1-O5-C5	3.02	116.23	112.19
5	C	1403	NAG	C1-O5-C5	3.01	116.21	112.19
5	B	1406	NAG	C3-C4-C5	2.98	115.64	110.23
5	A	1401	NAG	C1-O5-C5	2.90	116.08	112.19
5	A	1402	NAG	O5-C1-C2	-2.76	107.02	111.29
5	B	1404	NAG	C1-O5-C5	2.67	115.77	112.19
5	A	1402	NAG	C2-N2-C7	-2.63	119.37	122.90
5	B	1403	NAG	C1-O5-C5	2.48	115.51	112.19
5	B	1401	NAG	C2-N2-C7	2.47	126.21	122.90
5	B	1402	NAG	C1-O5-C5	2.36	115.35	112.19
5	B	1406	NAG	C1-O5-C5	2.32	115.29	112.19
5	A	1402	NAG	O4-C4-C3	-2.25	105.08	110.38
5	C	1408	NAG	O4-C4-C3	-2.18	105.25	110.38
5	A	1403	NAG	C1-O5-C5	2.11	115.02	112.19
5	B	1401	NAG	C4-C3-C2	-2.02	108.05	111.02
5	A	1410	NAG	C4-C3-C2	-2.00	108.08	111.02

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1401	NAG	C1-C2-N2-C7
5	C	1401	NAG	O5-C5-C6-O6
5	C	1401	NAG	C4-C5-C6-O6
5	A	1403	NAG	O5-C5-C6-O6
5	A	1402	NAG	C8-C7-N2-C2
5	B	1404	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	B	1402	NAG	O5-C5-C6-O6
5	A	1406	NAG	C4-C5-C6-O6
5	A	1406	NAG	O5-C5-C6-O6
5	A	1403	NAG	C4-C5-C6-O6
5	A	1402	NAG	O7-C7-N2-C2
5	C	1403	NAG	O5-C5-C6-O6
5	B	1407	NAG	C4-C5-C6-O6
5	B	1404	NAG	C3-C2-N2-C7
5	C	1403	NAG	C3-C2-N2-C7
5	B	1402	NAG	C4-C5-C6-O6

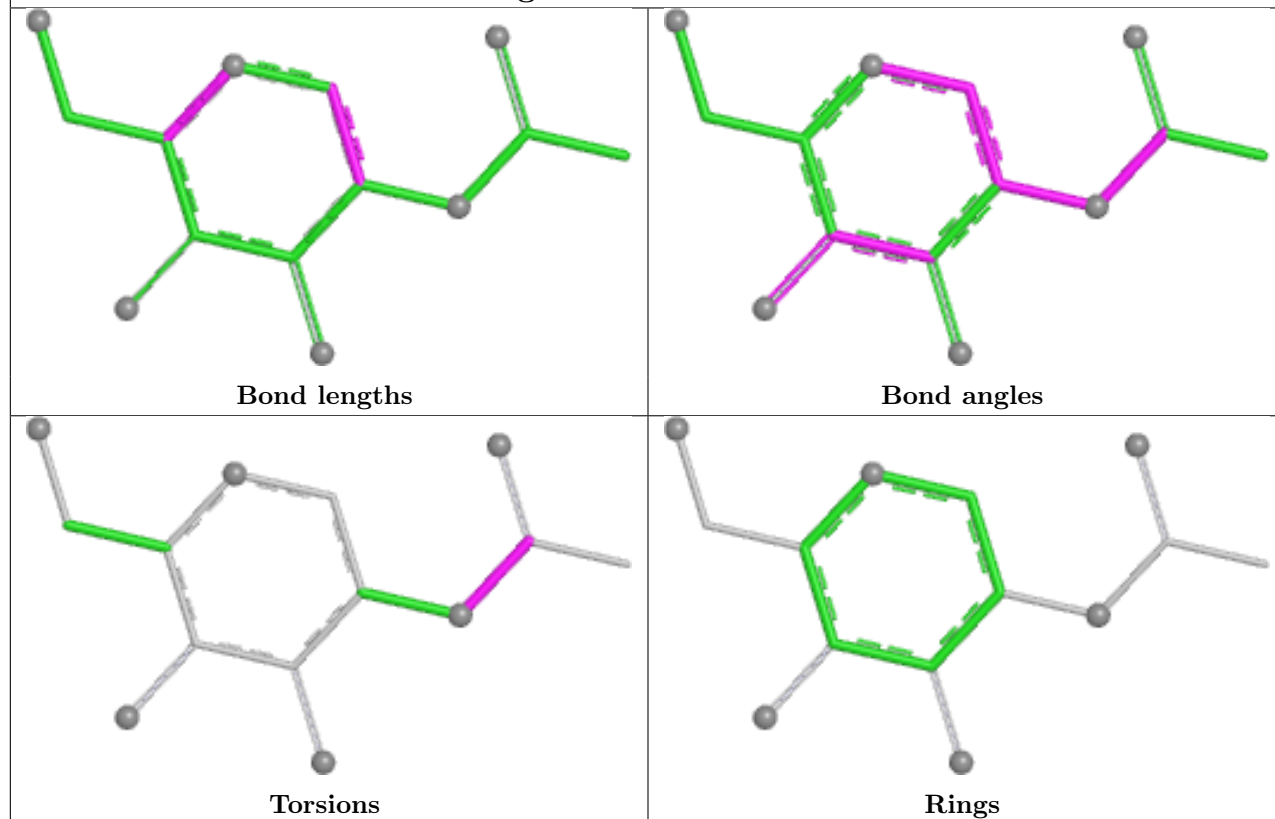
There are no ring outliers.

4 monomers are involved in 9 short contacts:

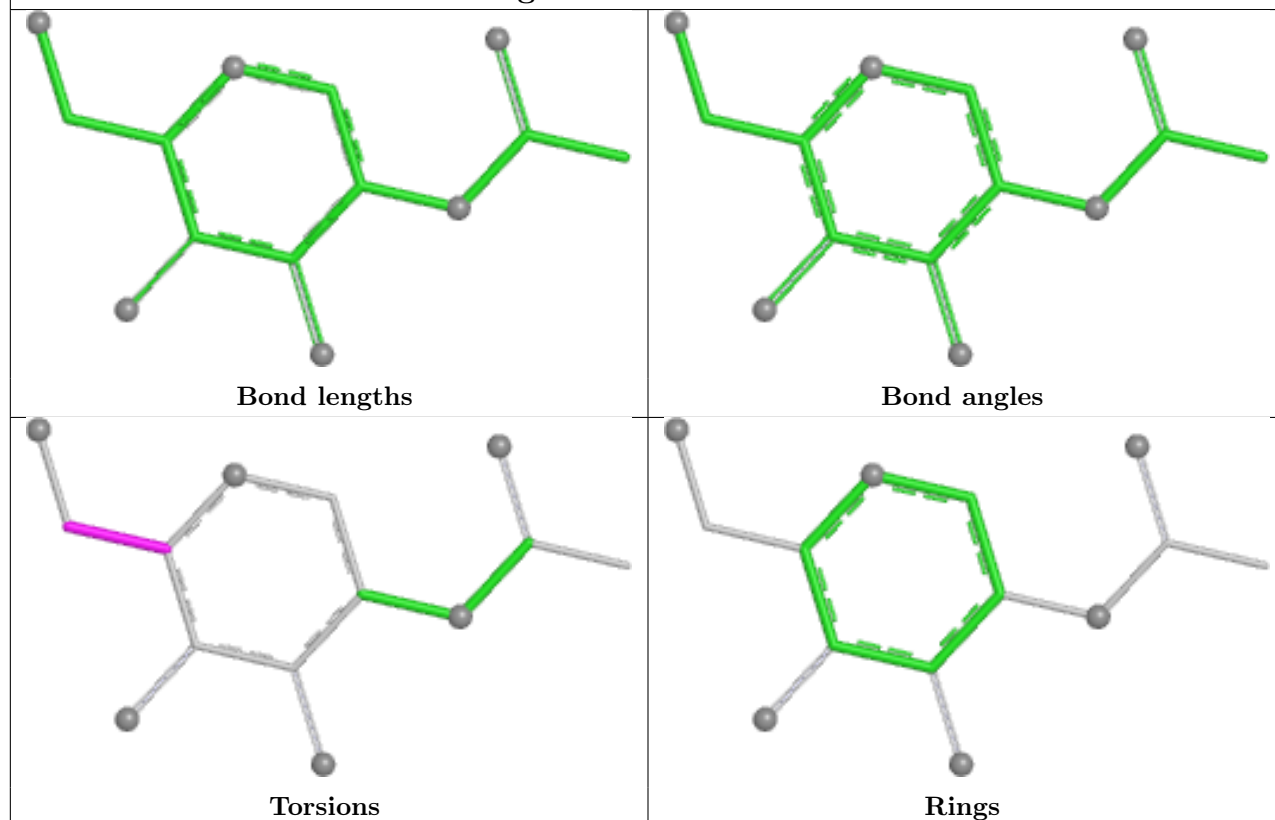
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1403	NAG	1	0
5	B	1401	NAG	1	0
5	B	1402	NAG	6	0
5	C	1407	NAG	1	0

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

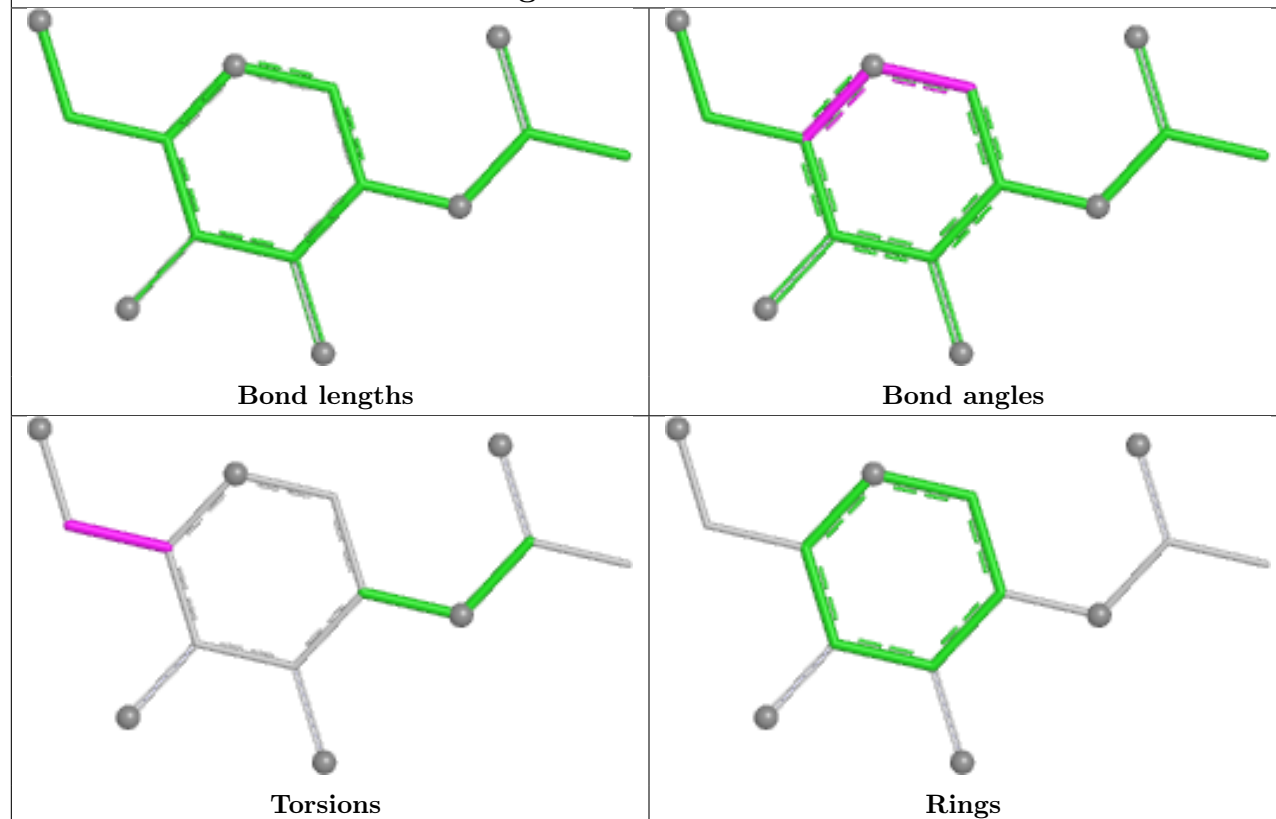
Ligand NAG A 1402



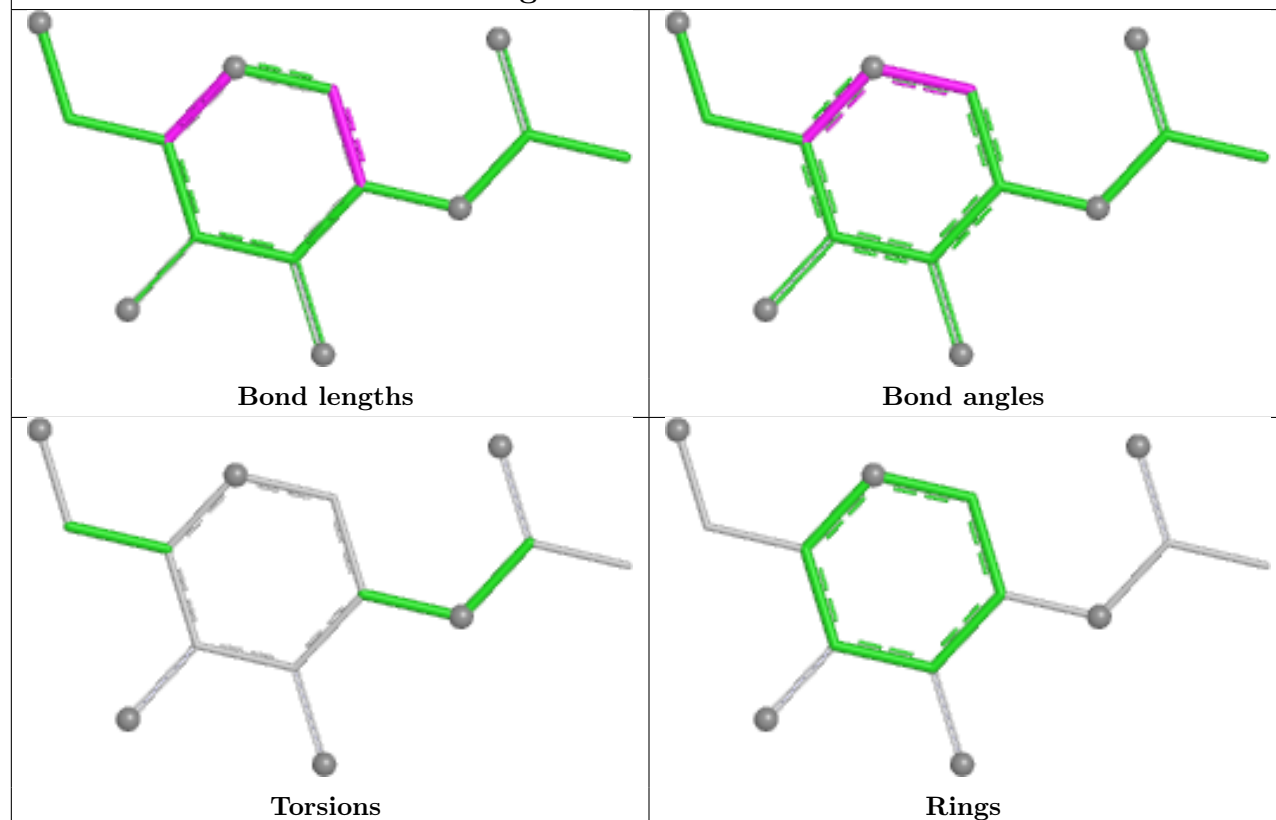
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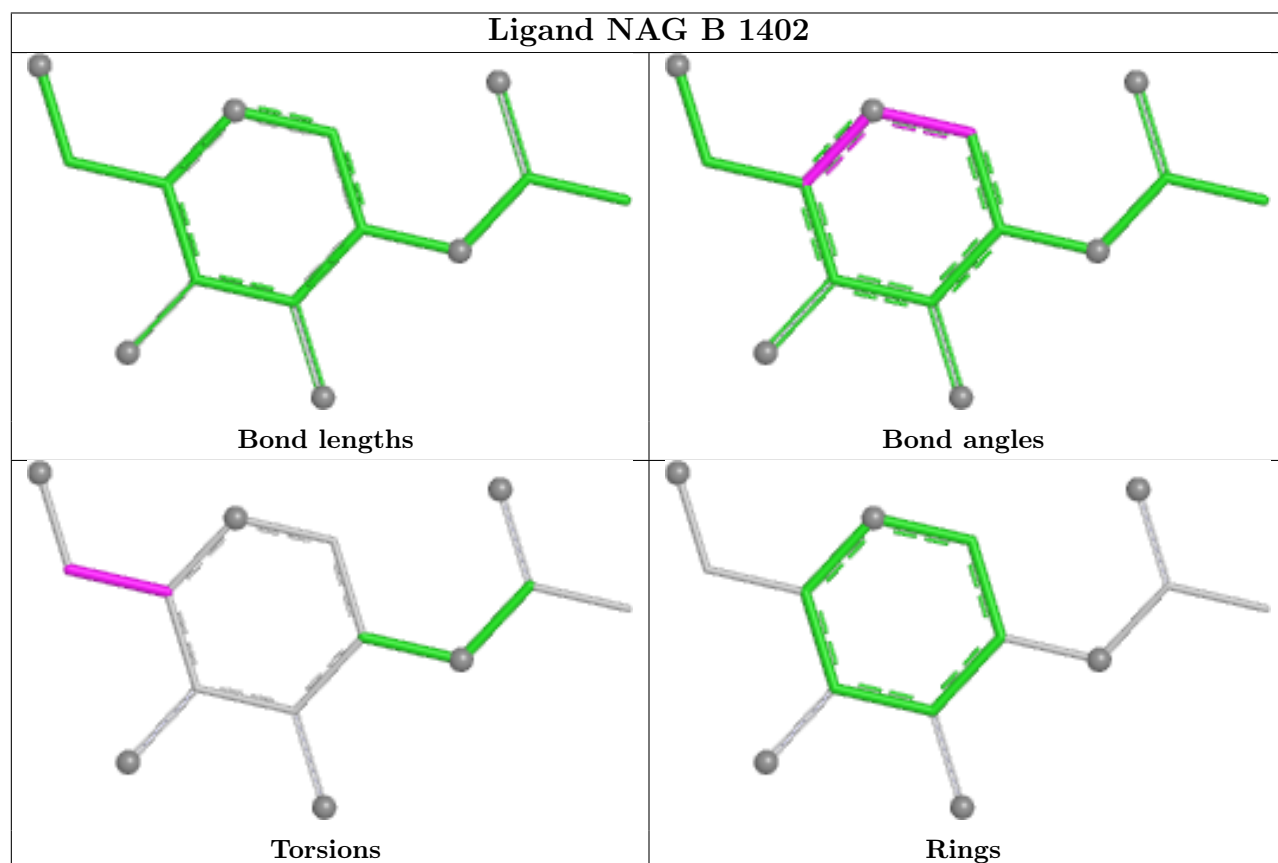
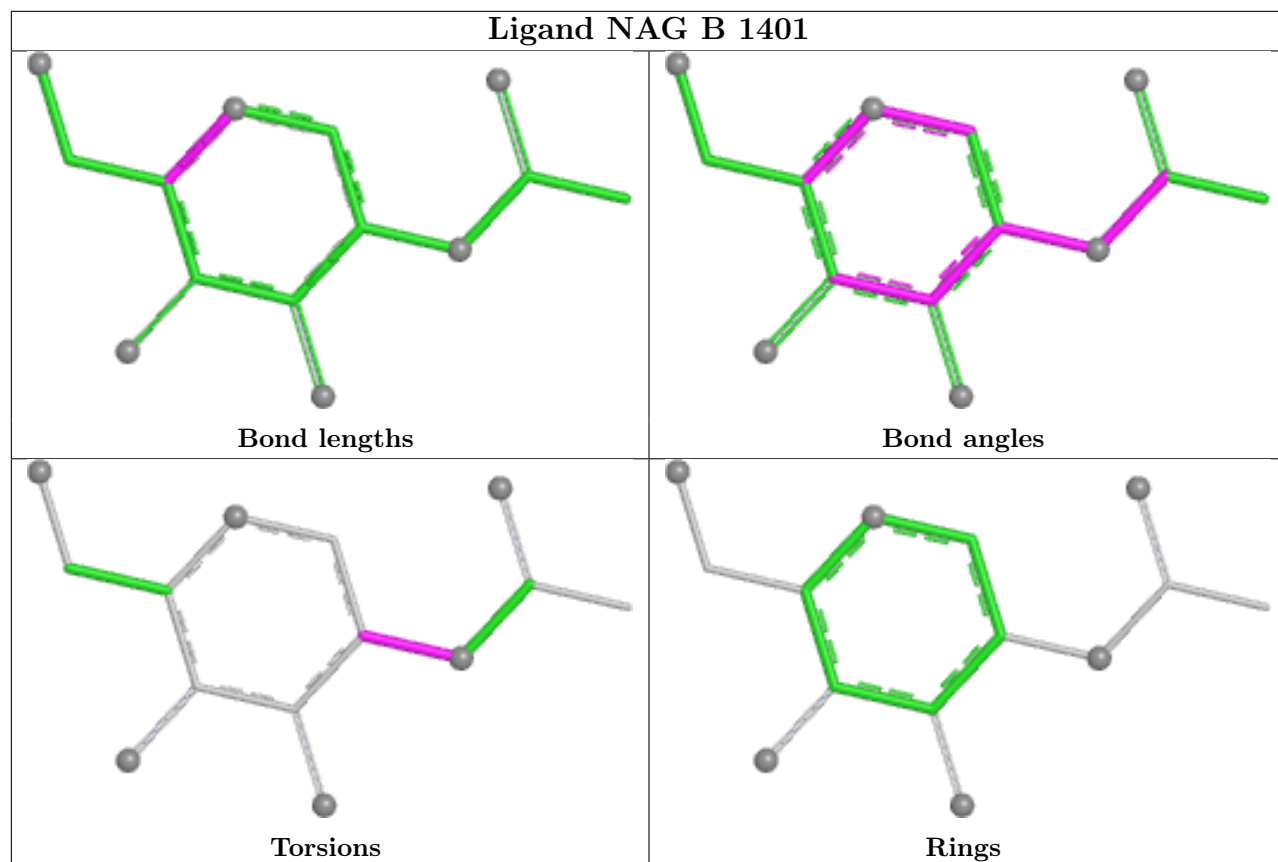


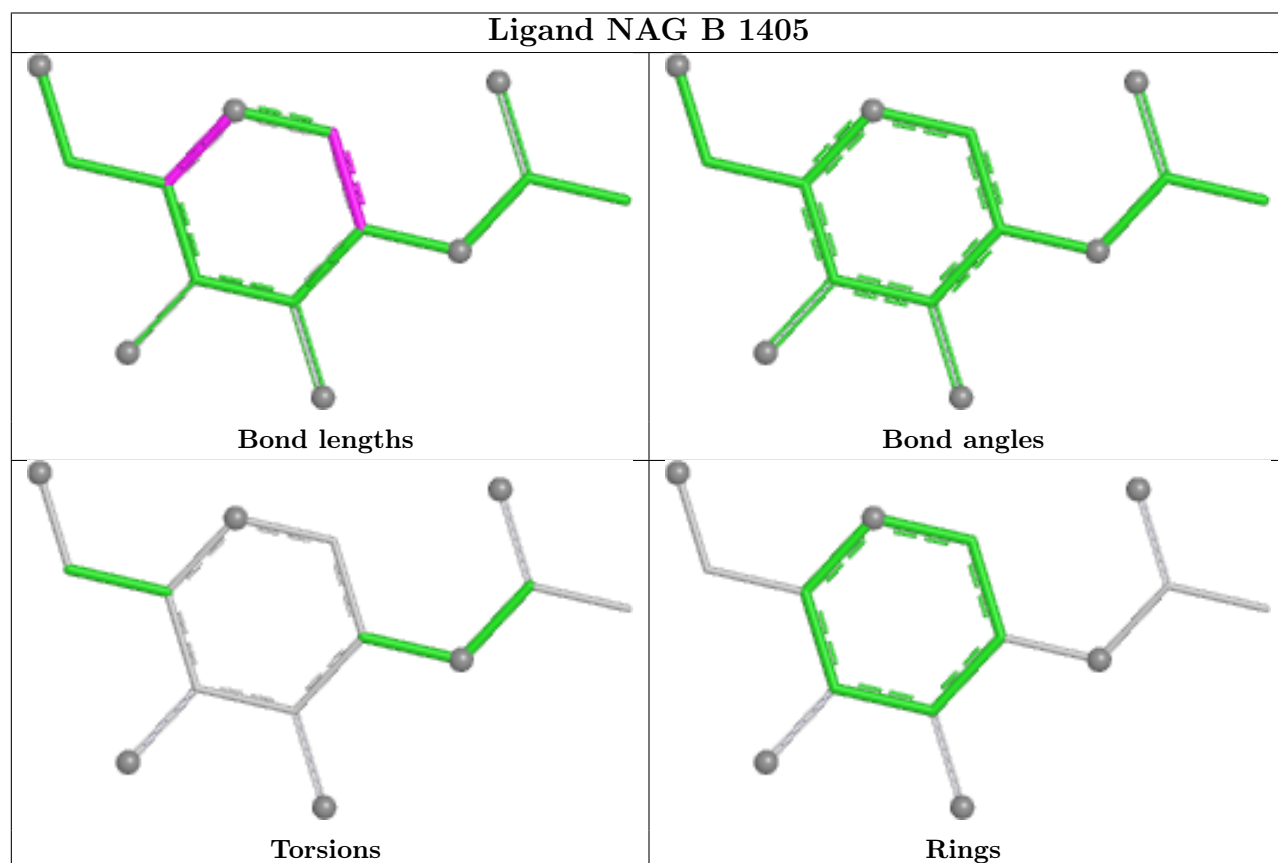
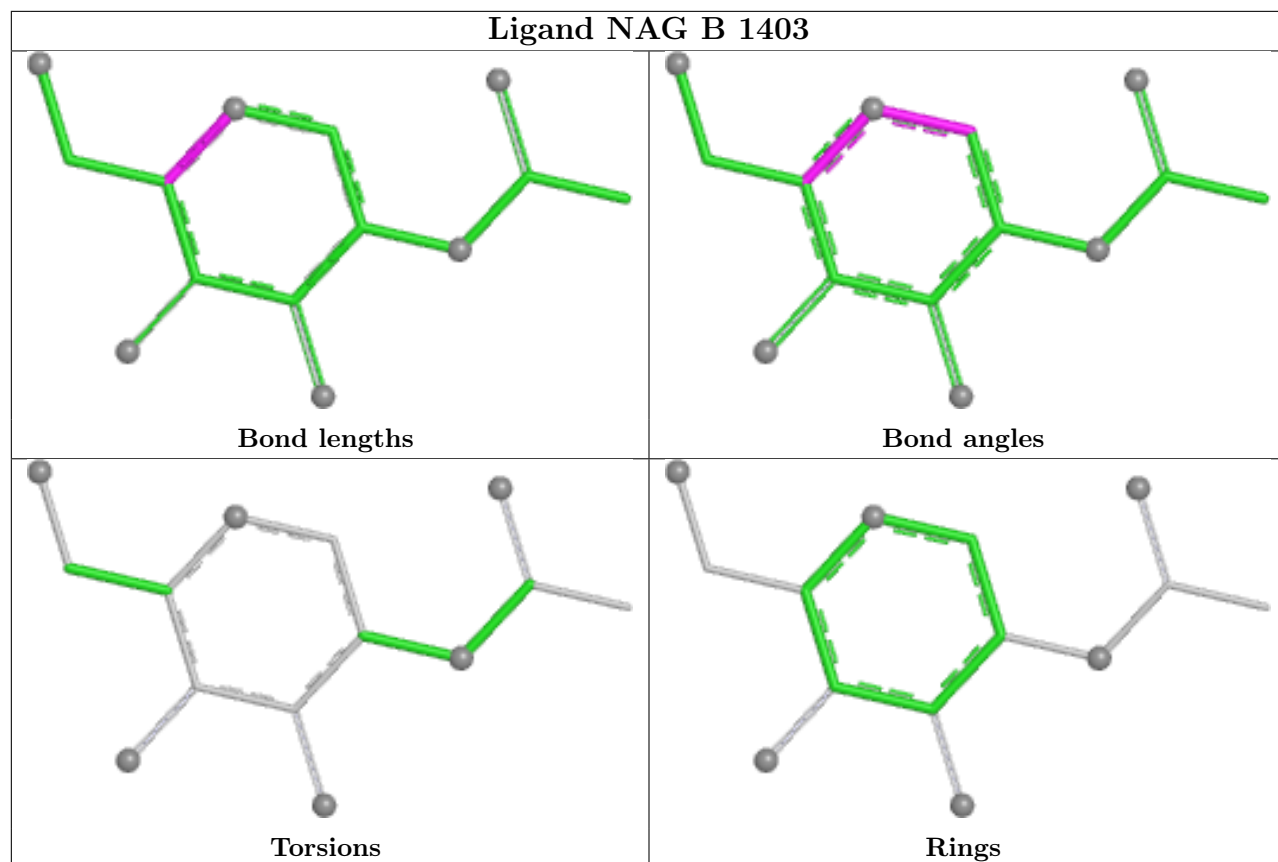
Ligand NAG A 1403

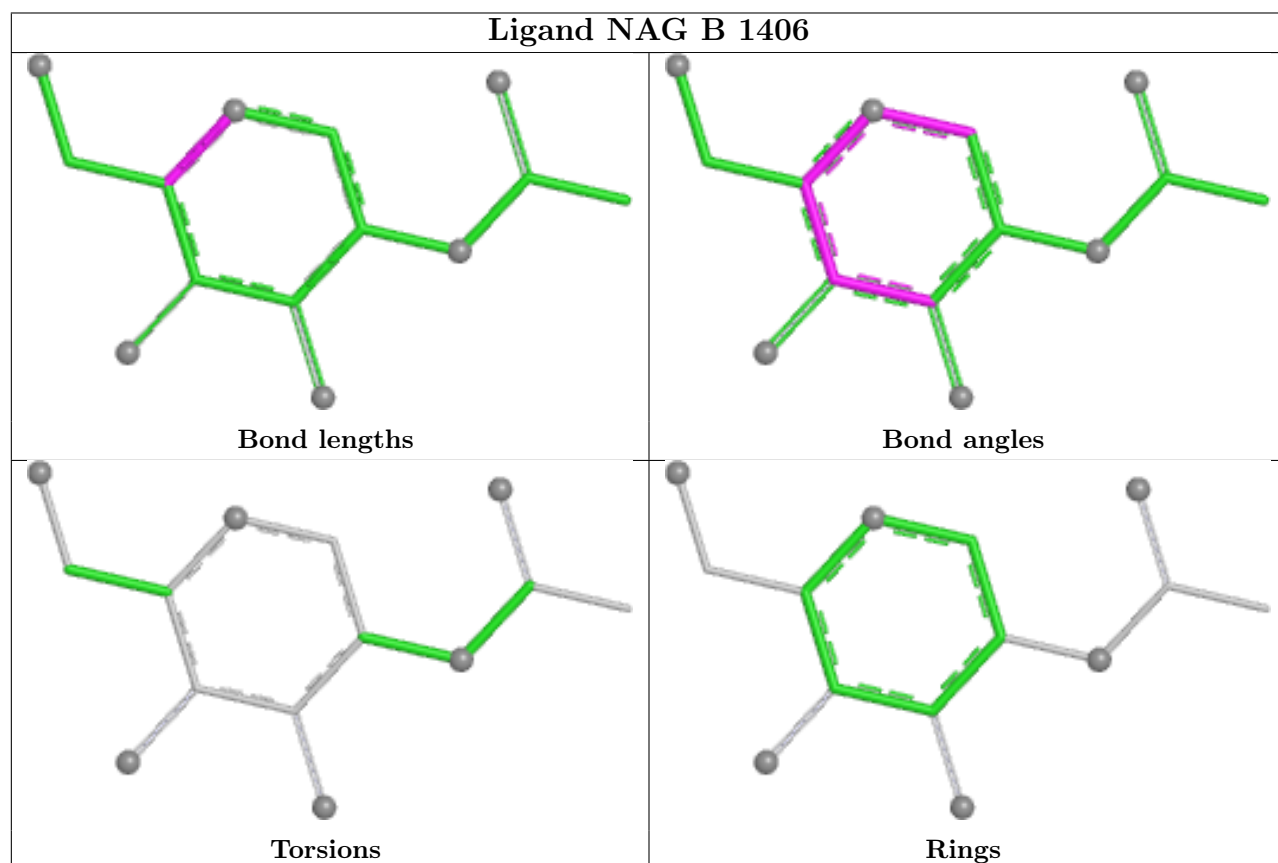
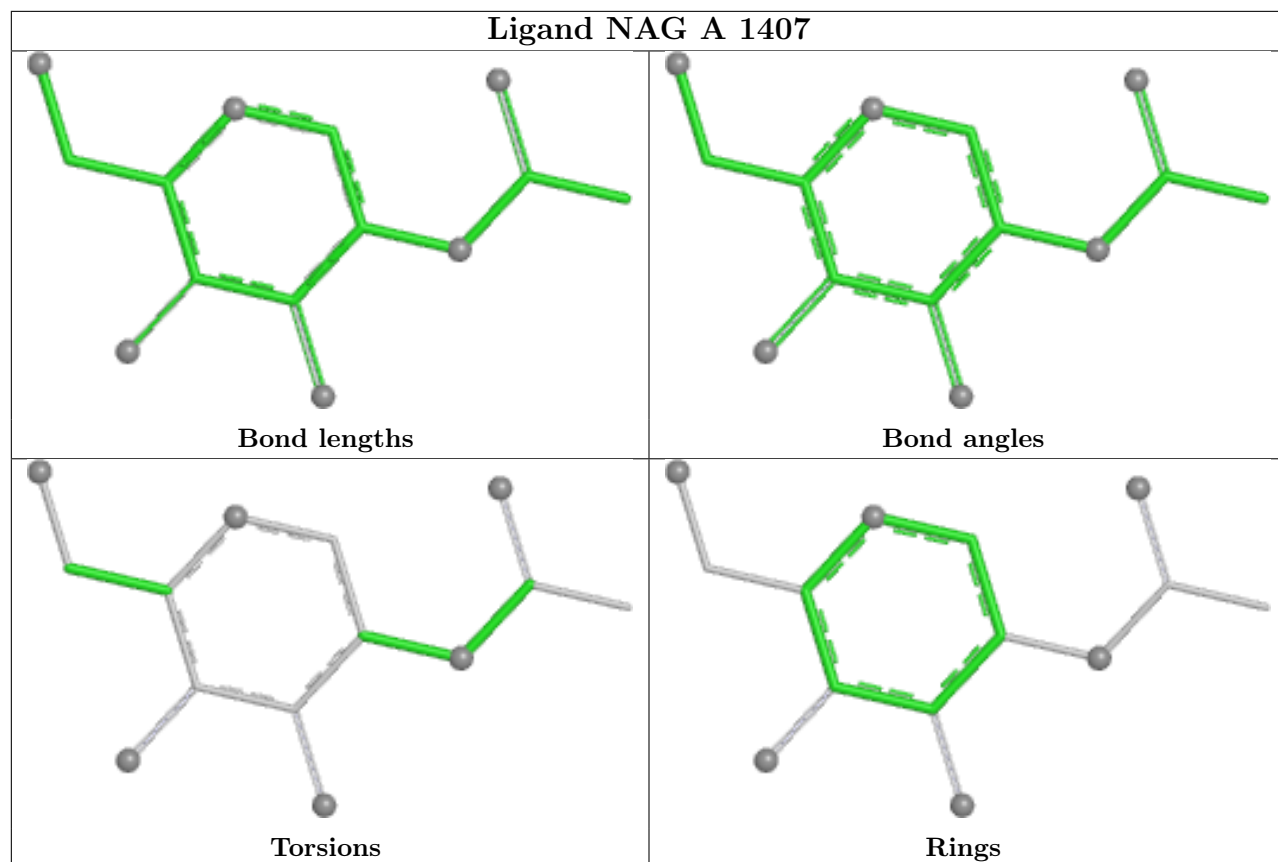


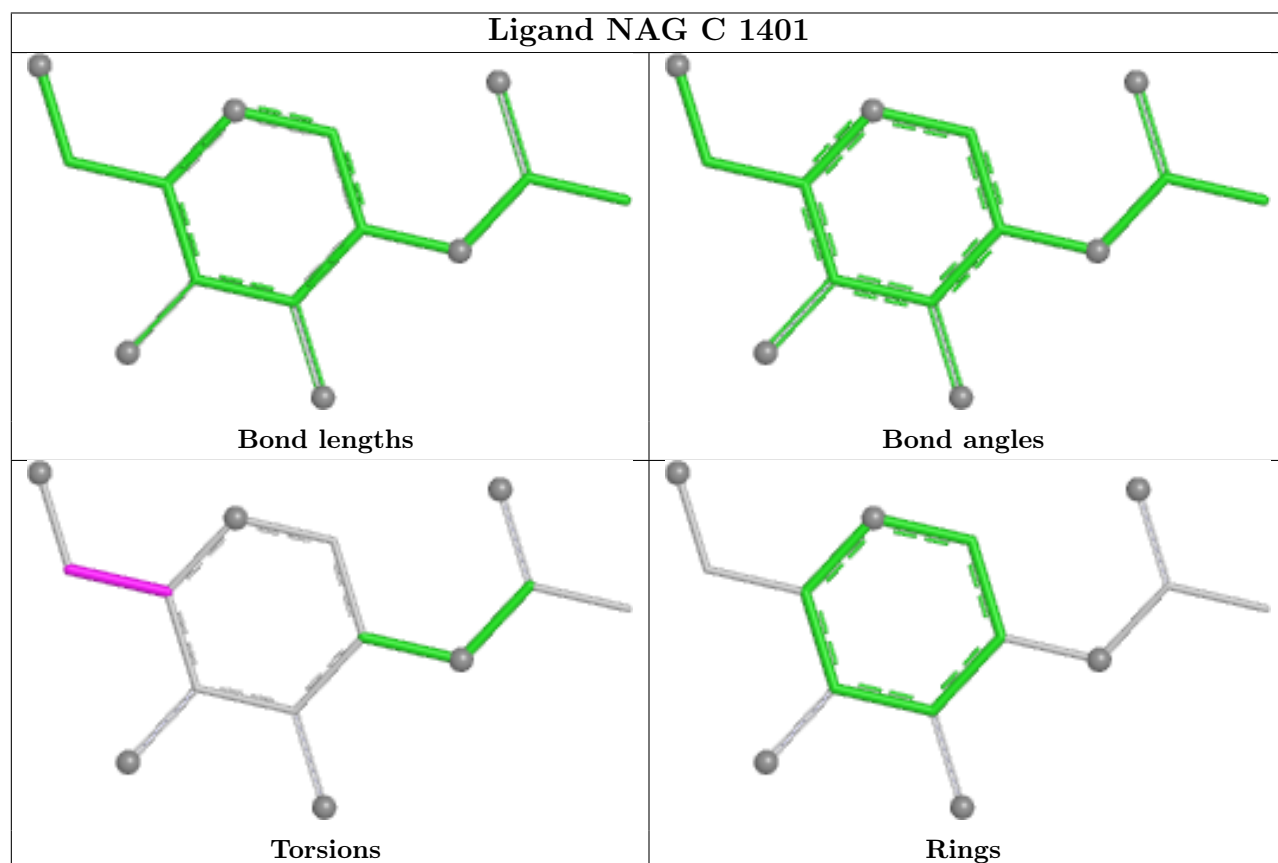
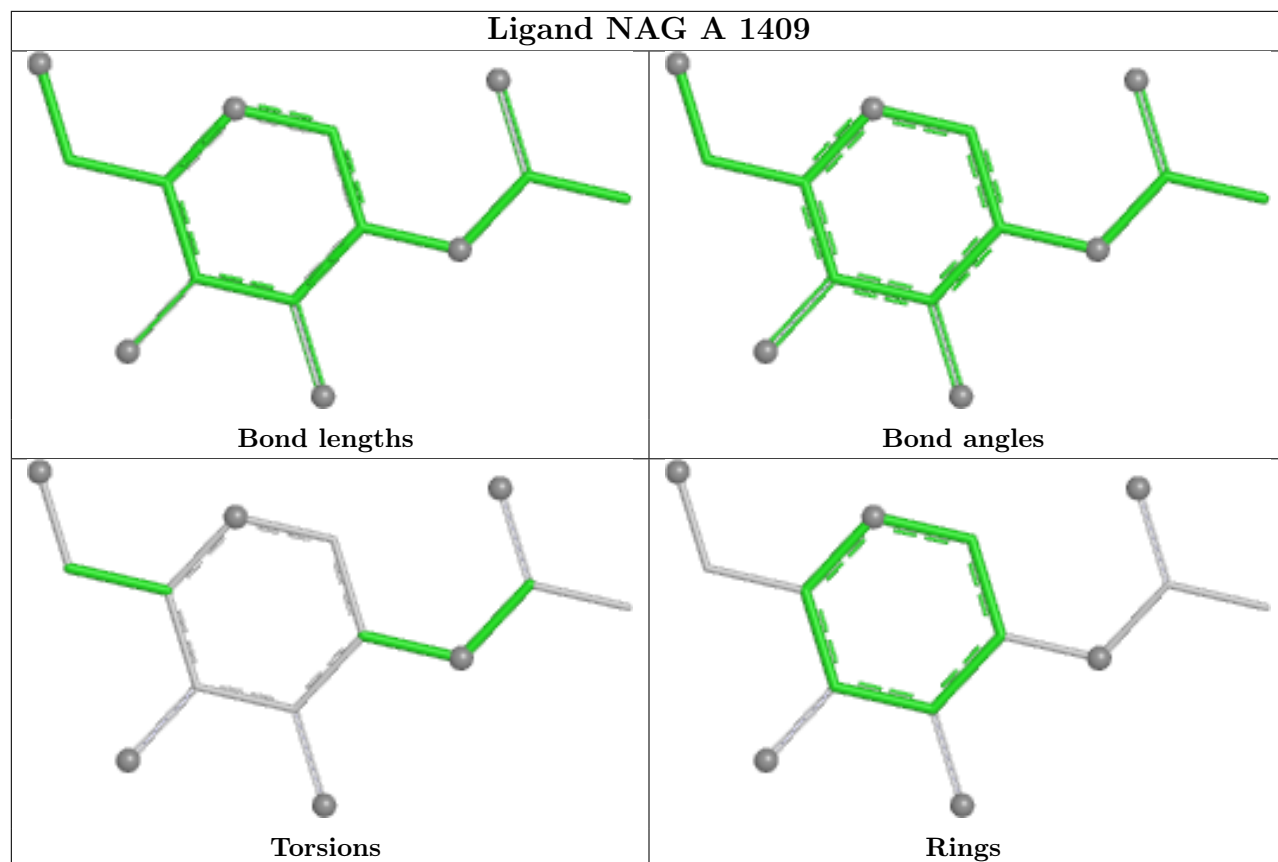
Ligand NAG B 1408



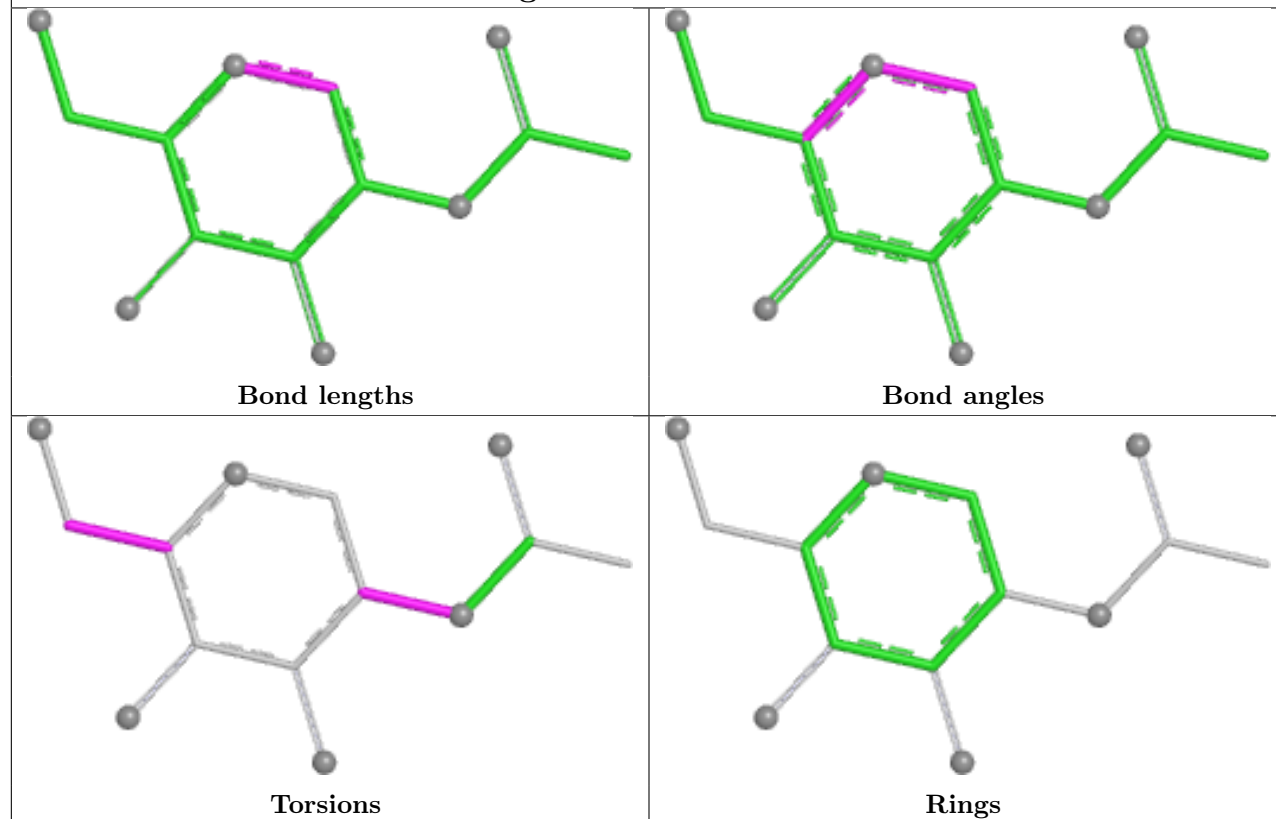




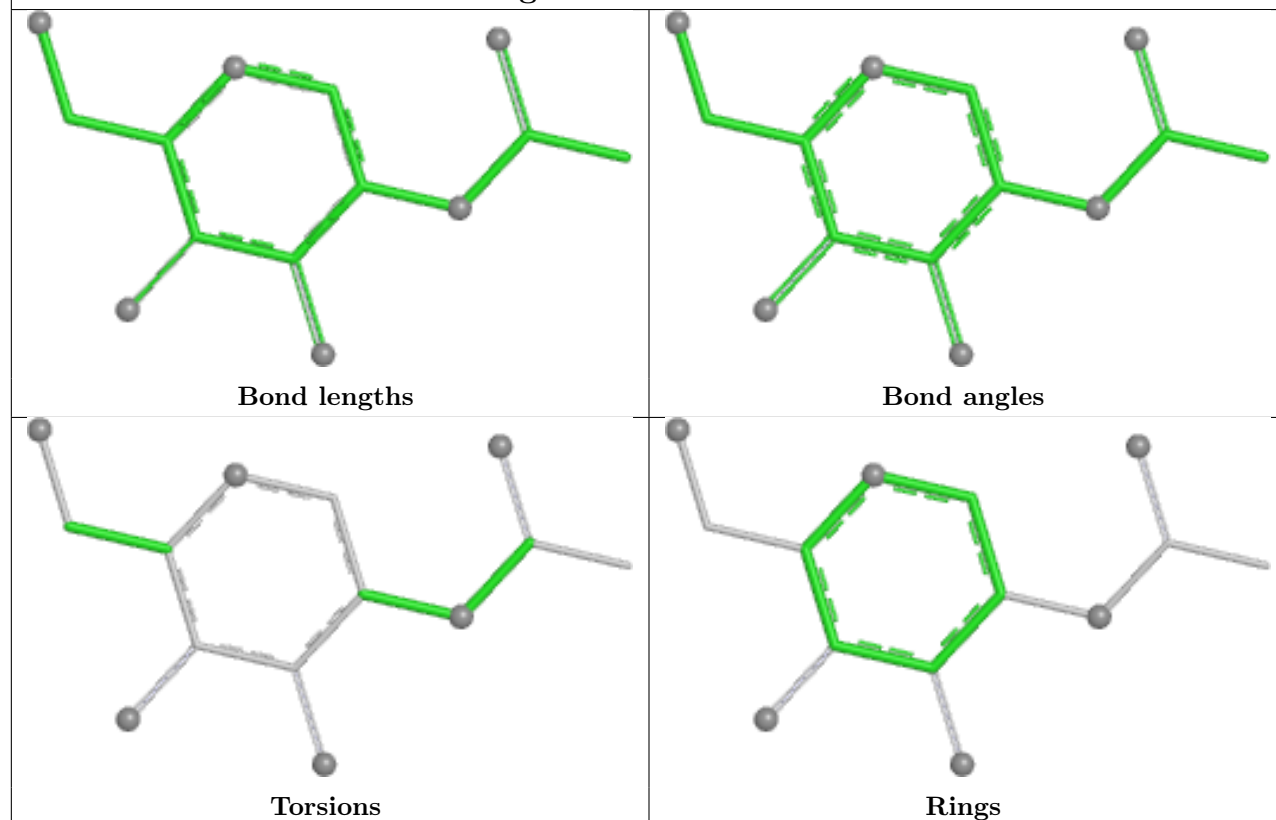


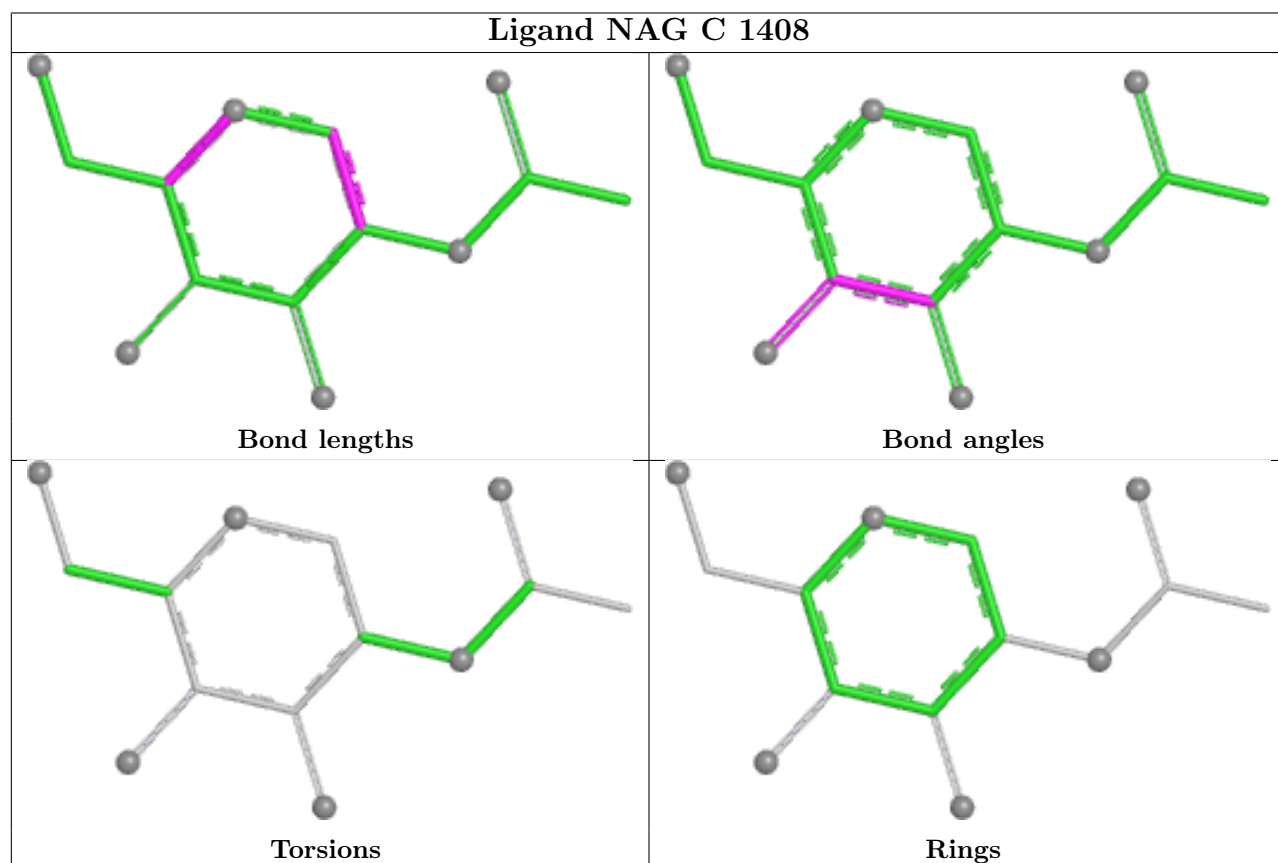
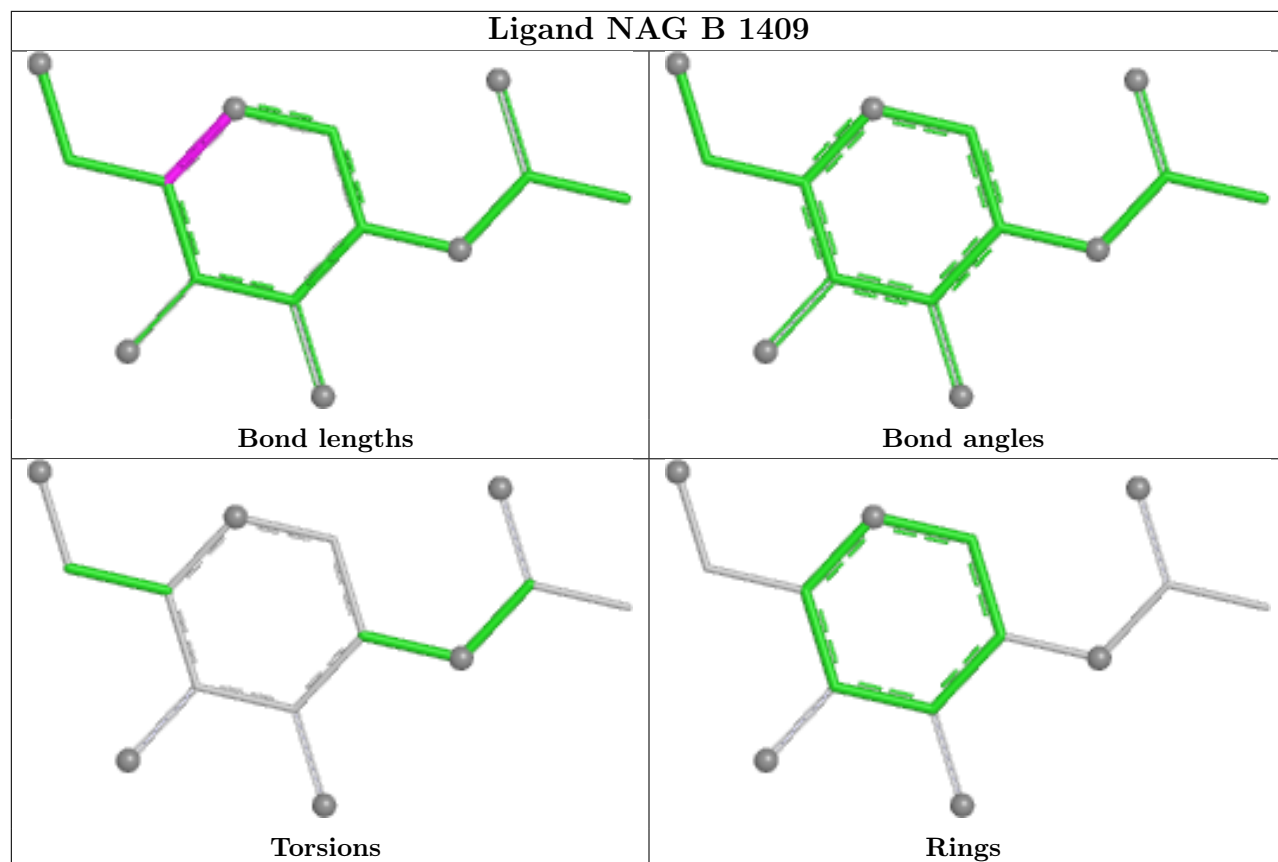


Ligand NAG C 1403

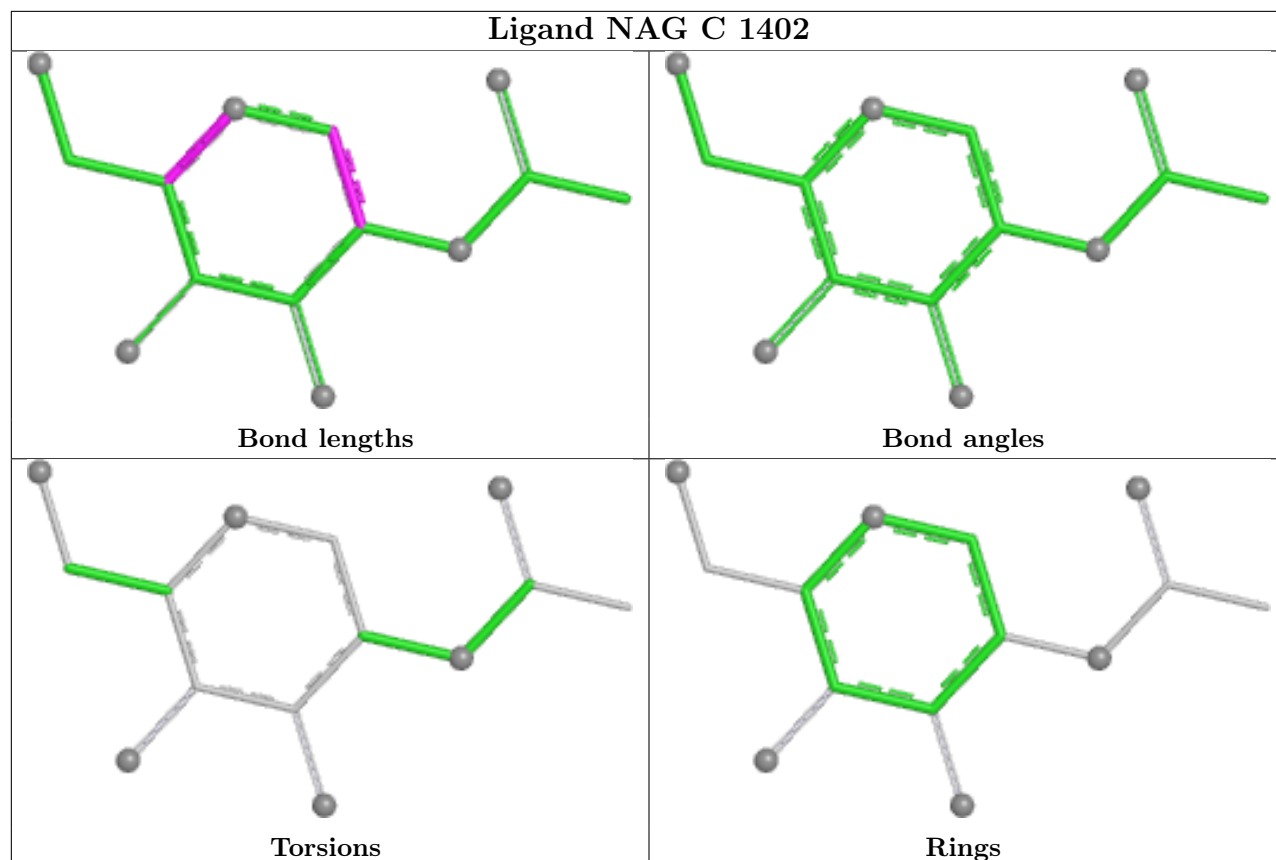


Ligand NAG C 1405

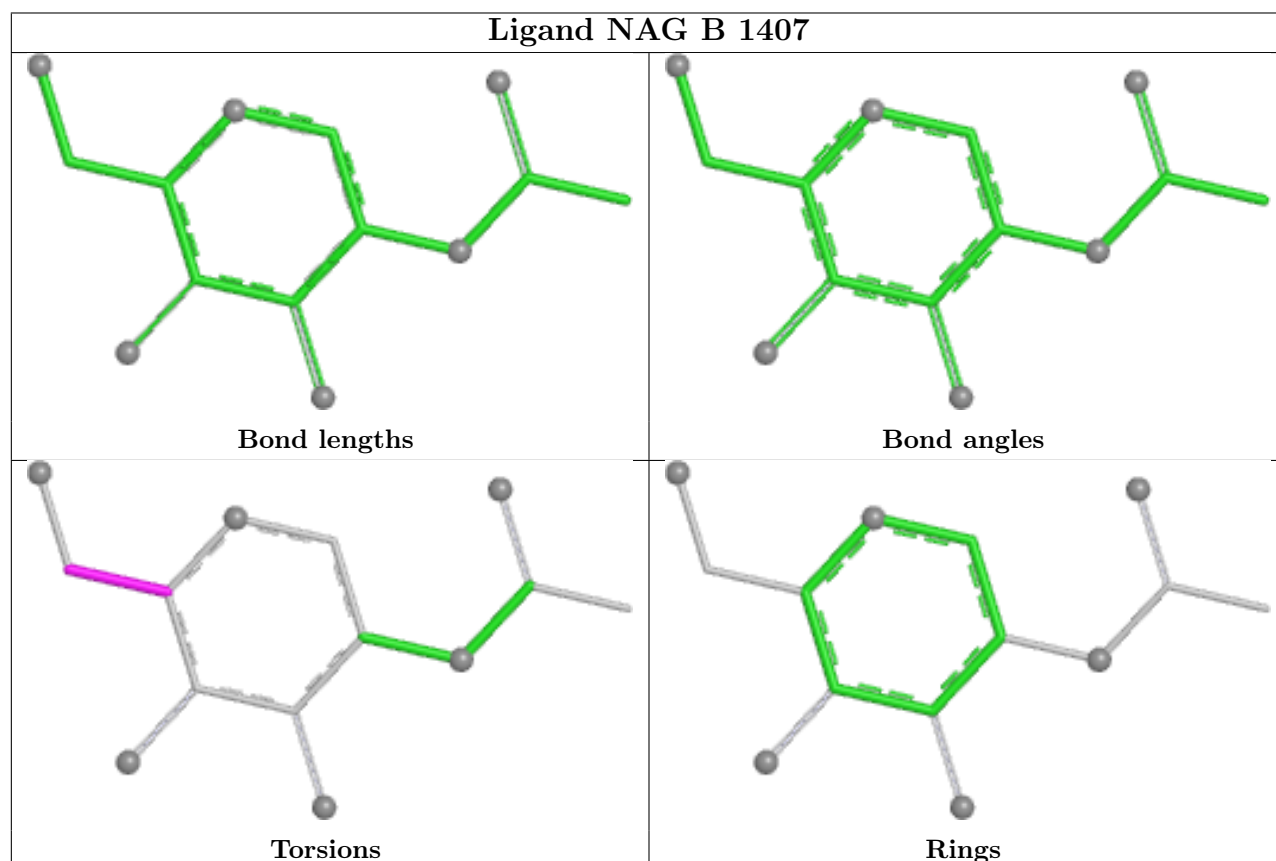




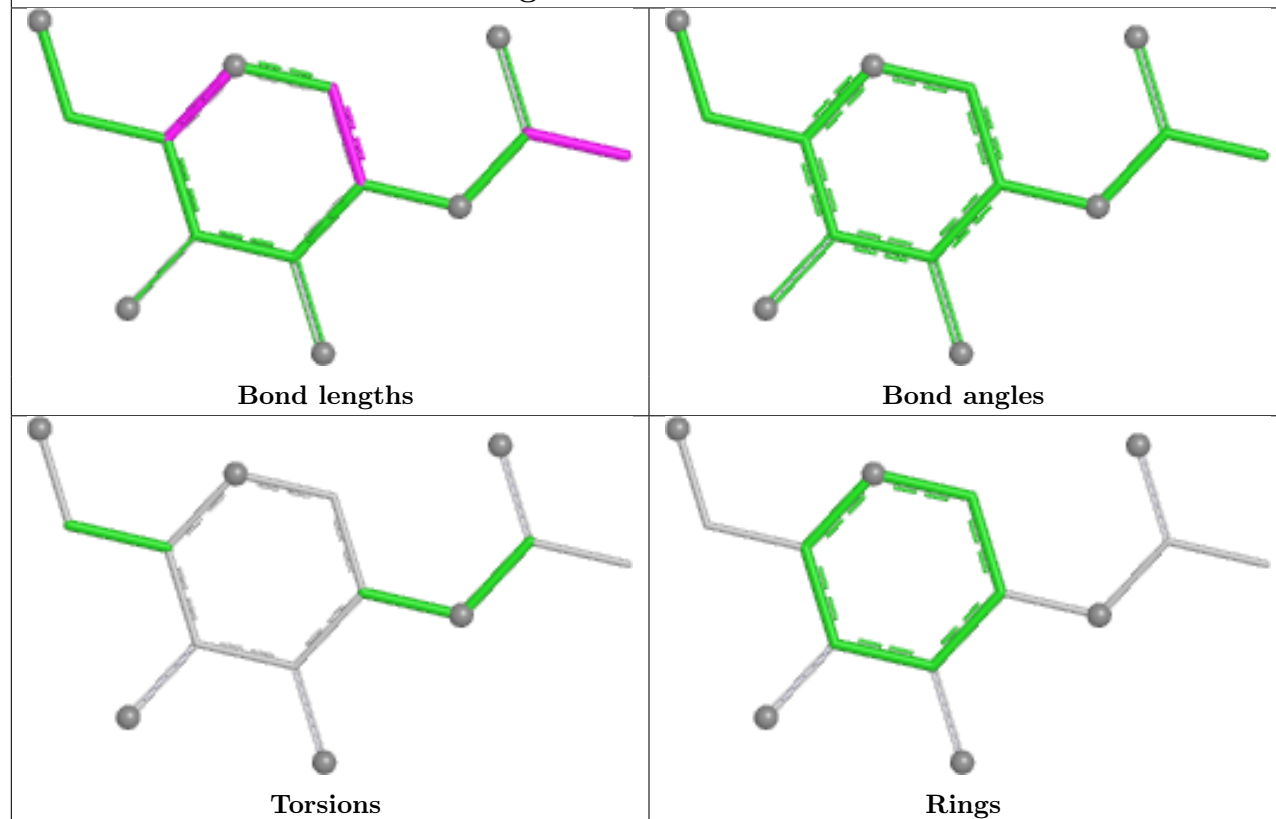
Ligand NAG C 1402



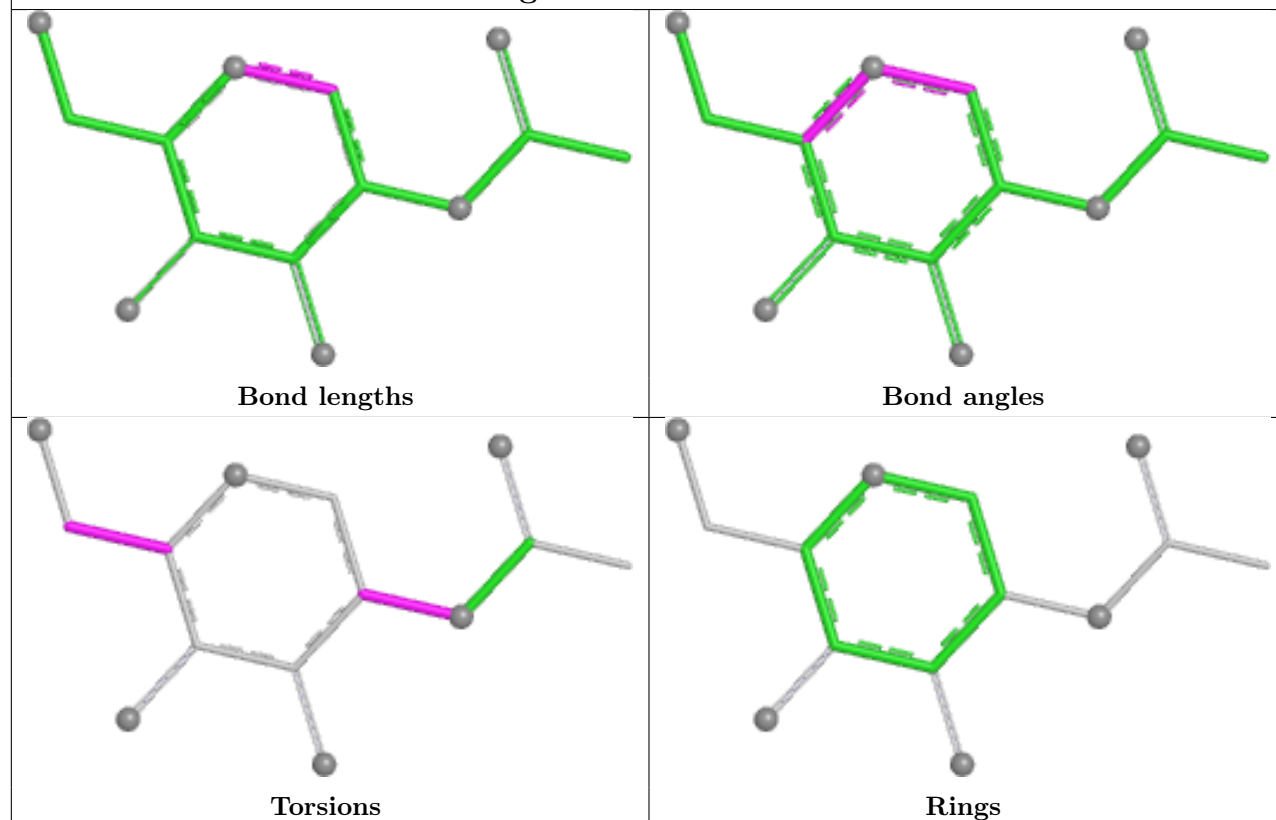
Ligand NAG B 1407



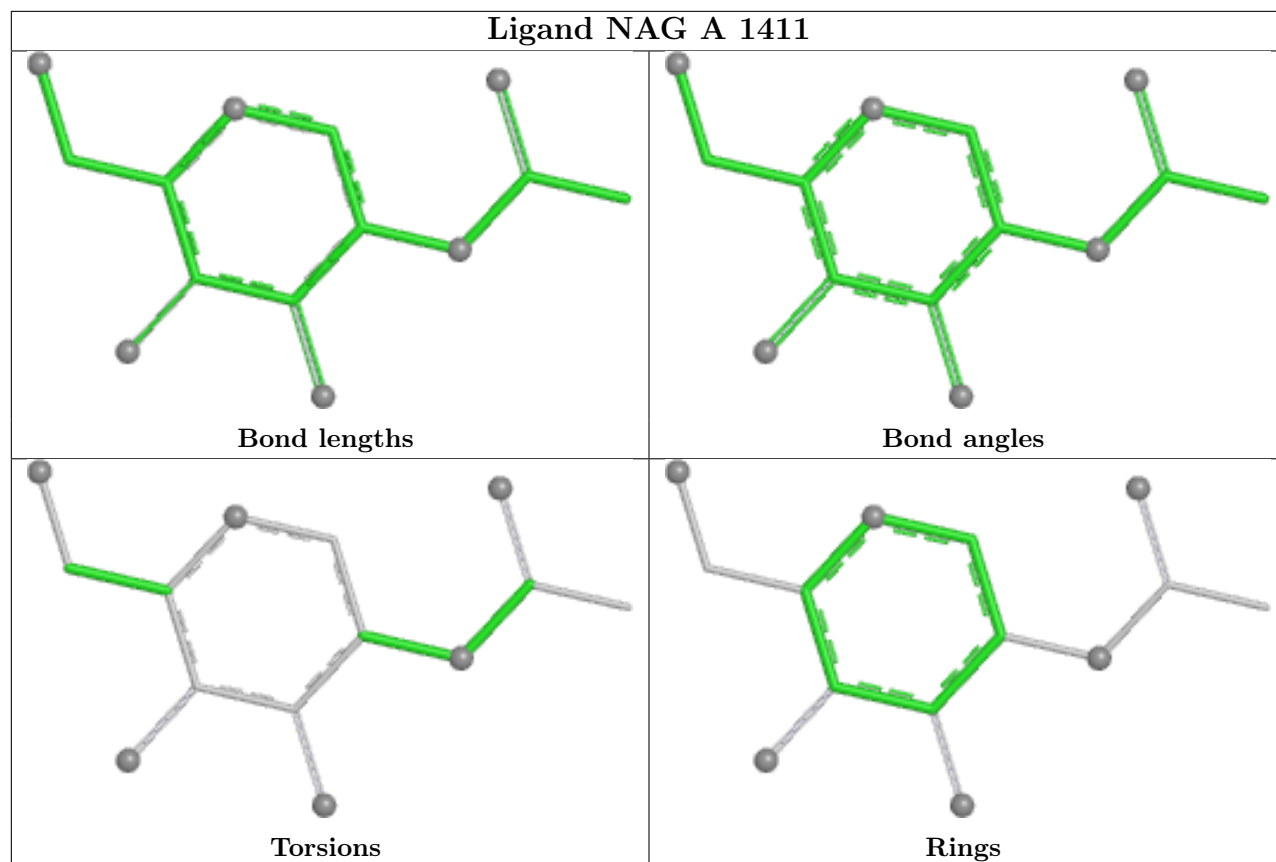
Ligand NAG C 1407



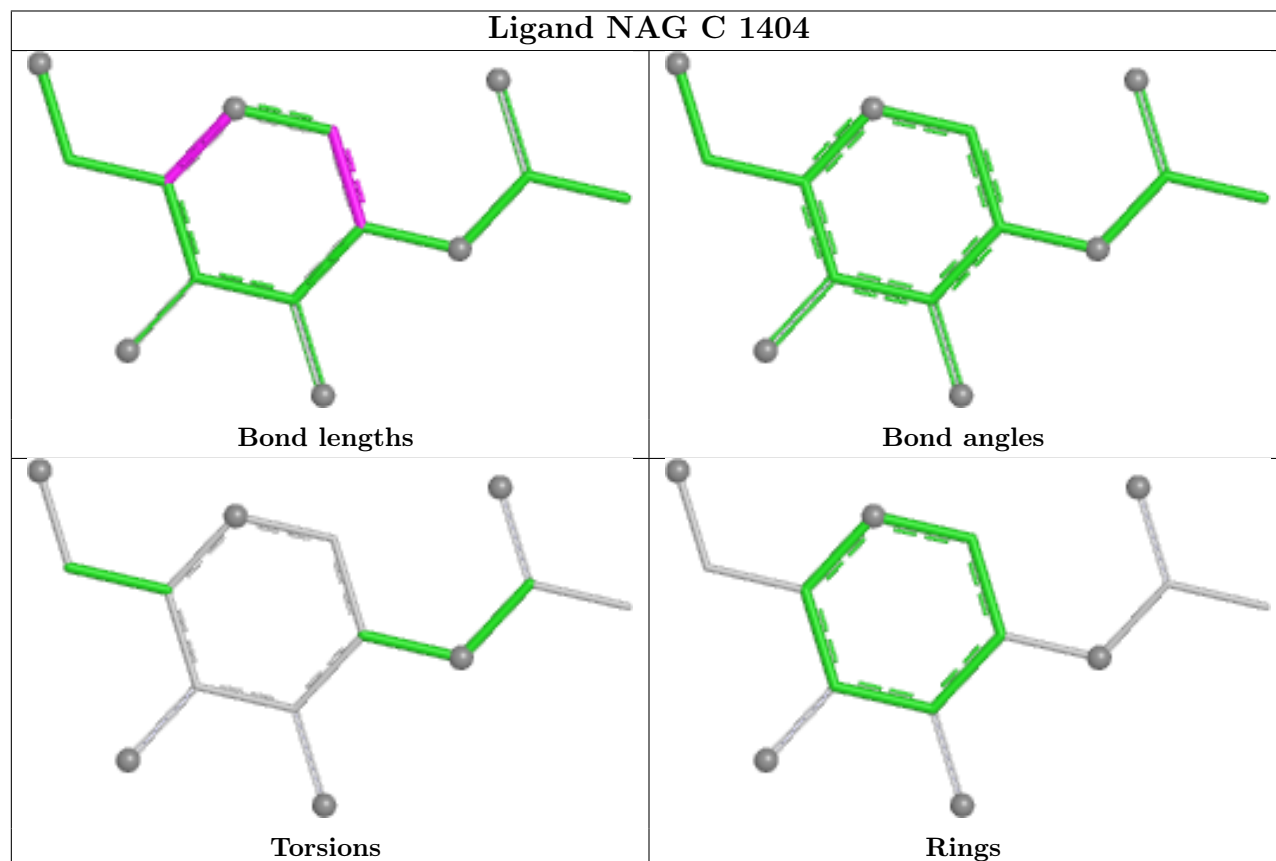
Ligand NAG B 1404

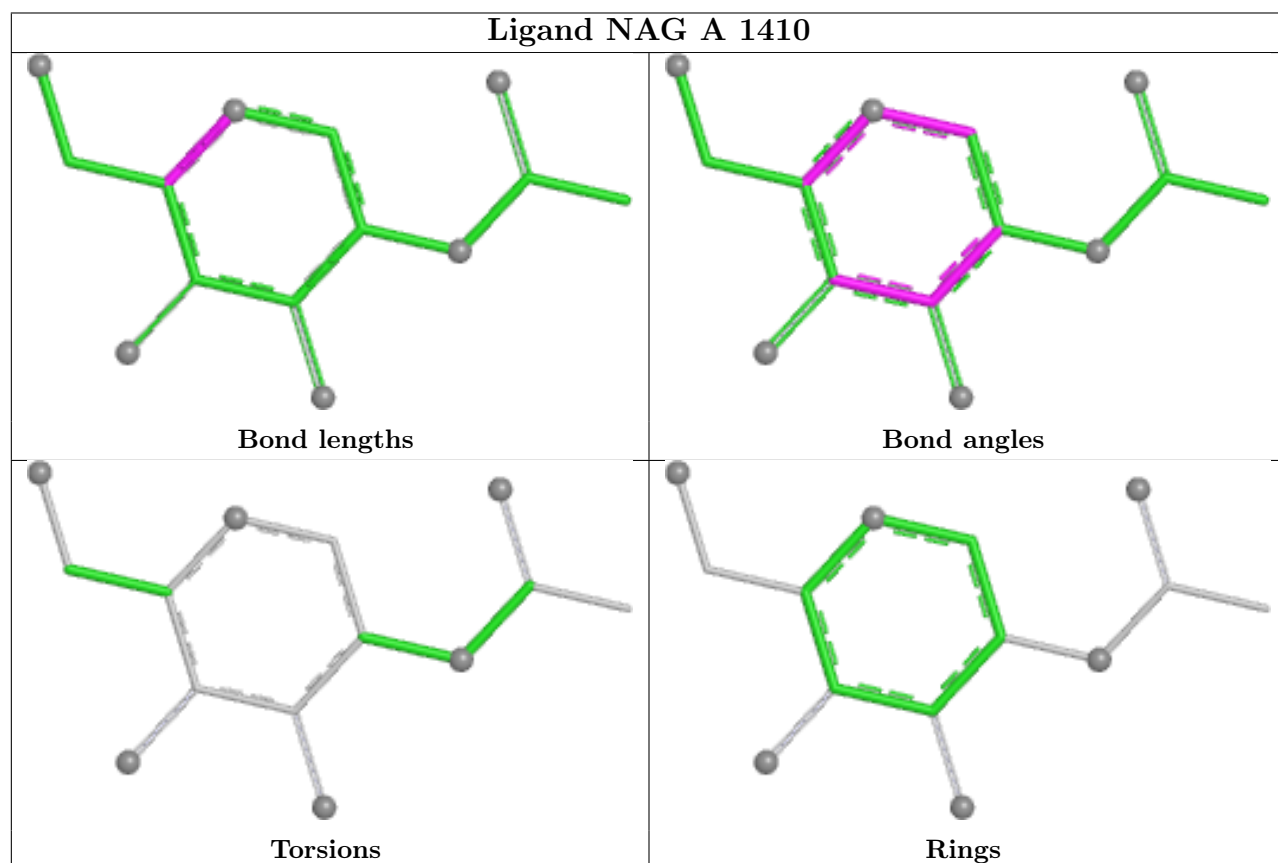
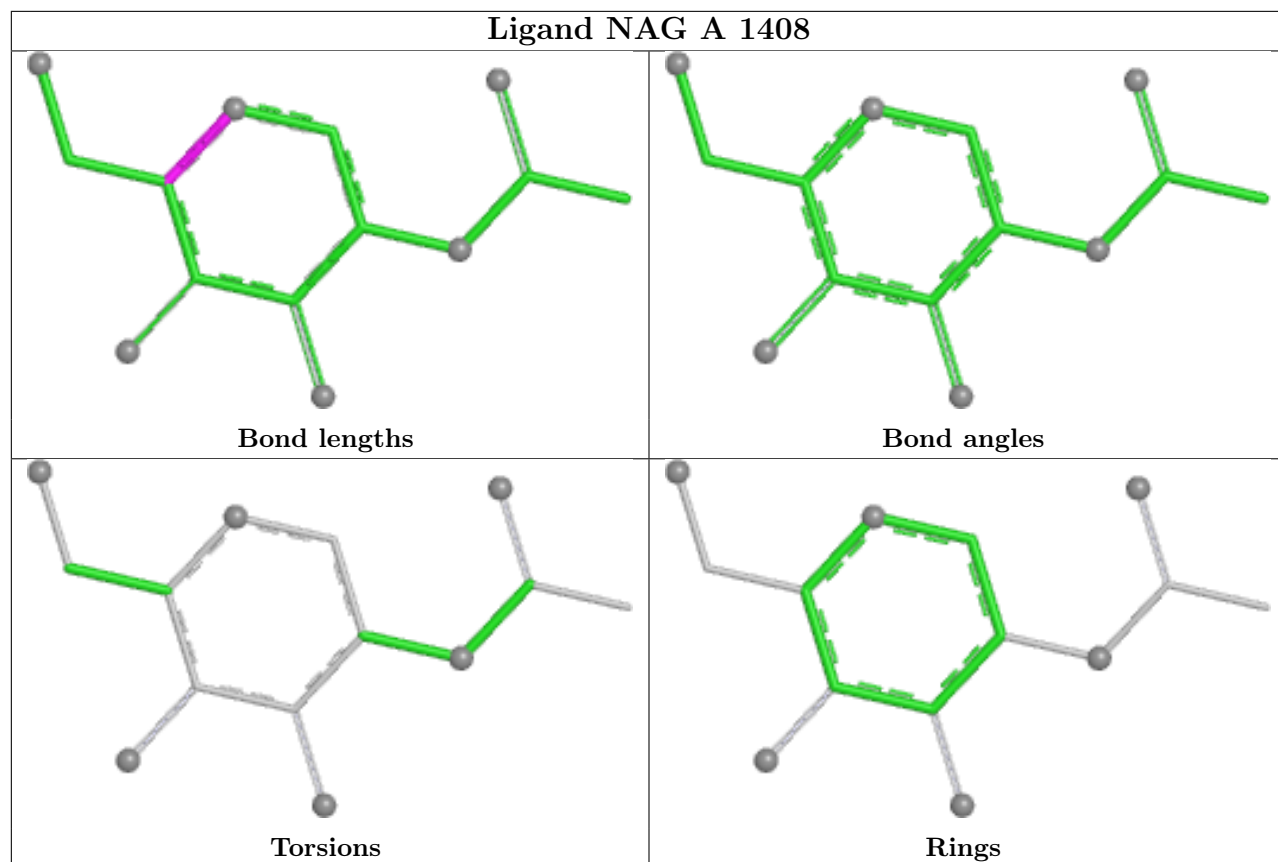


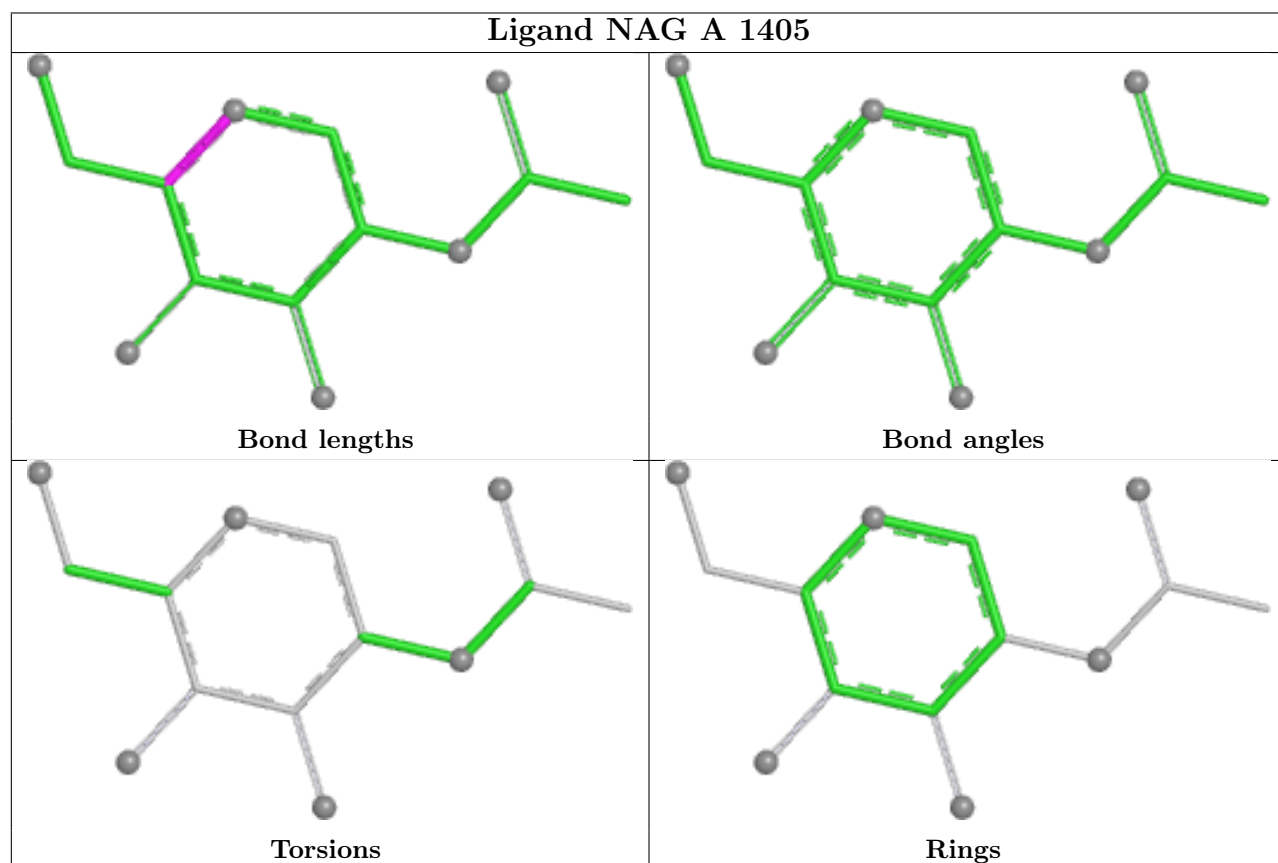
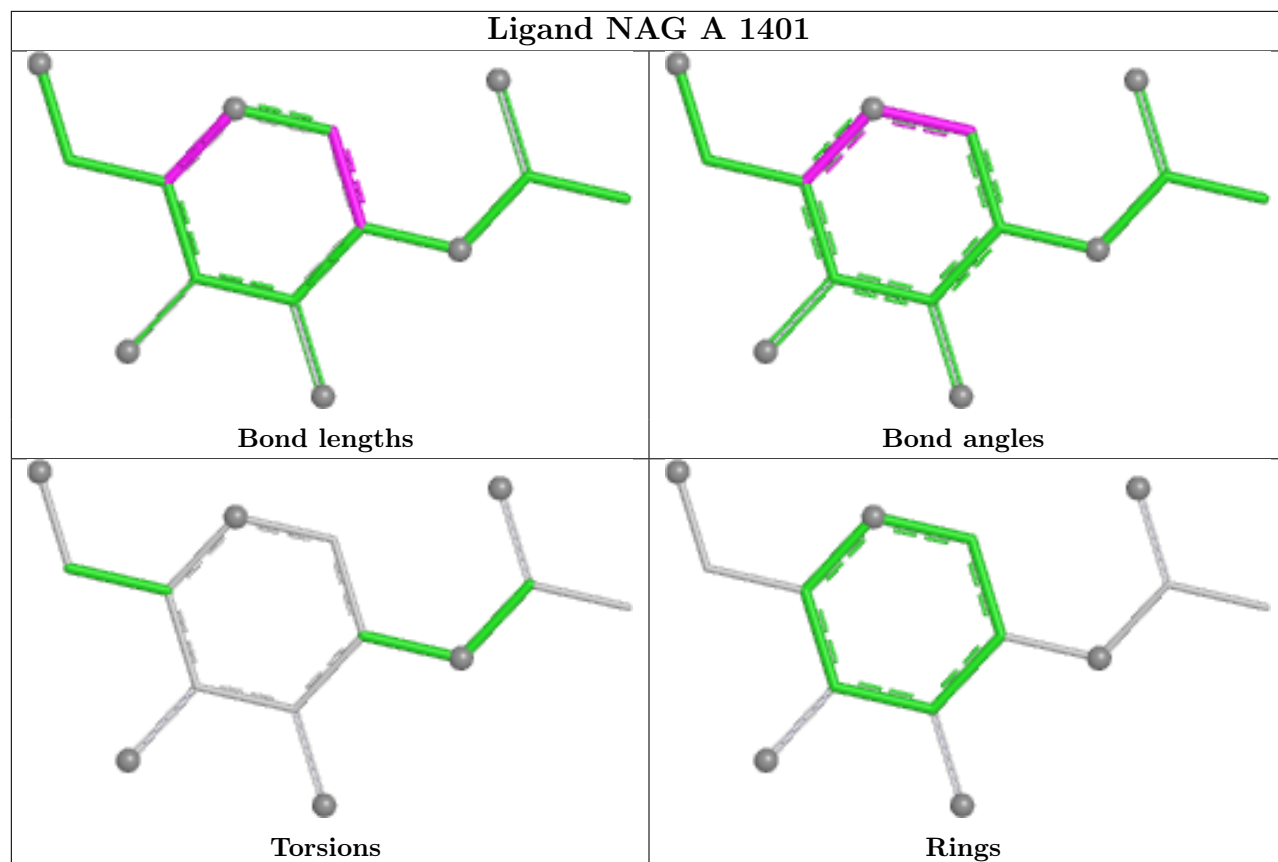
Ligand NAG A 1411



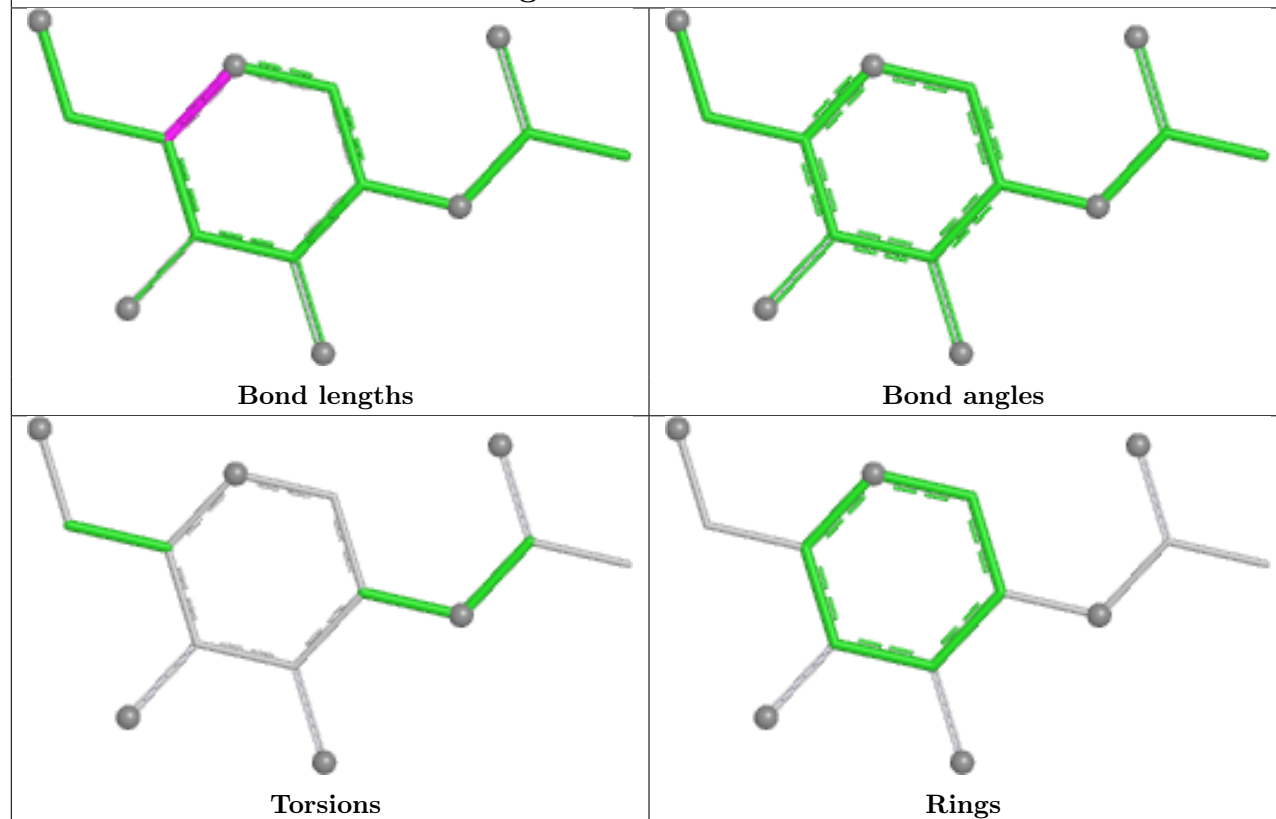
Ligand NAG C 1404



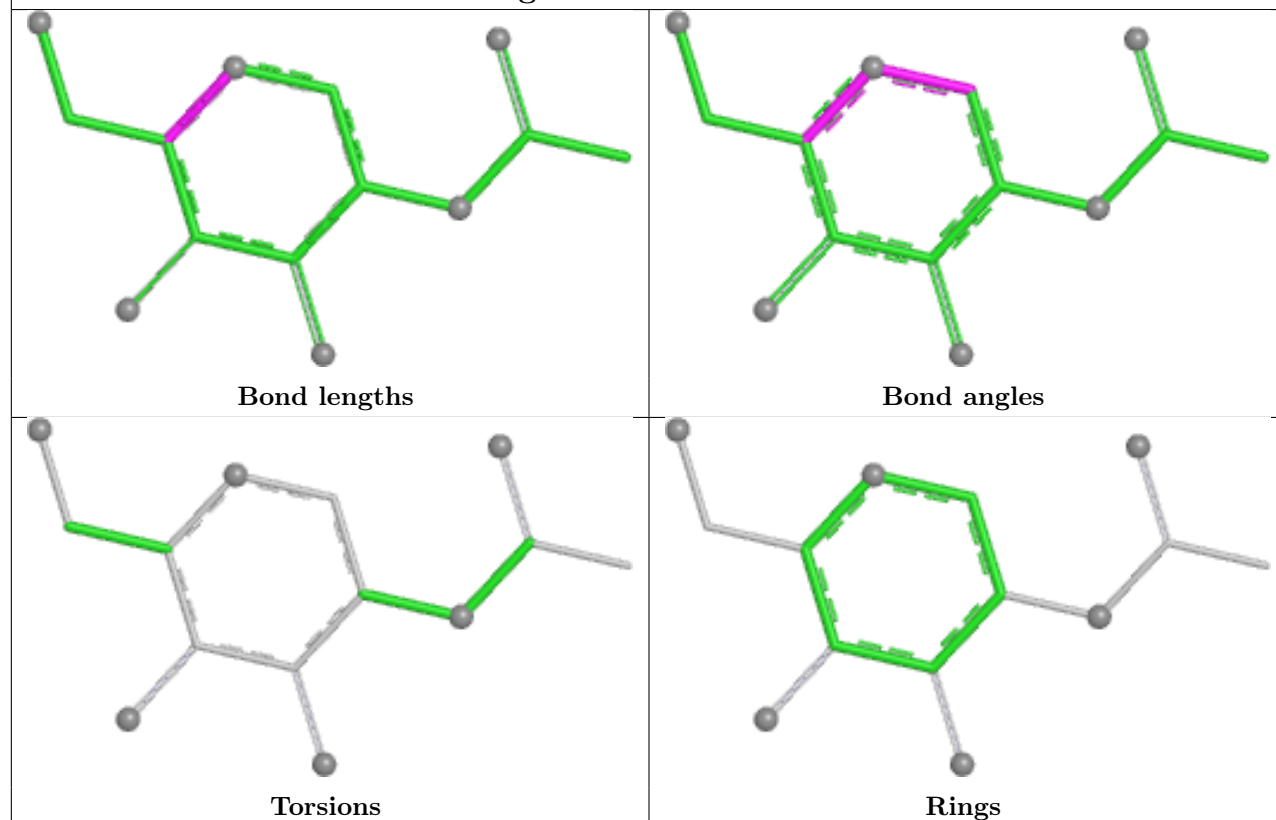




Ligand NAG C 1406



Ligand NAG A 1404



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

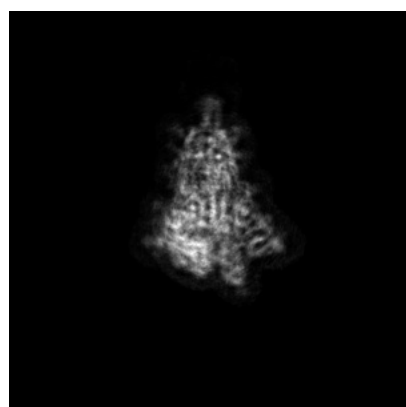
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24988. These allow visual inspection of the internal detail of the map and identification of artifacts.

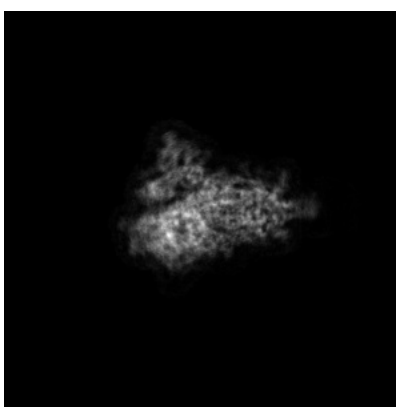
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

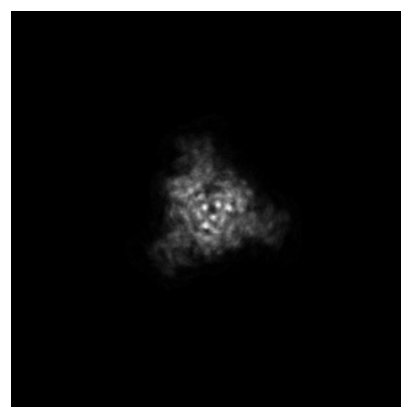
6.1.1 Primary map



X



Y

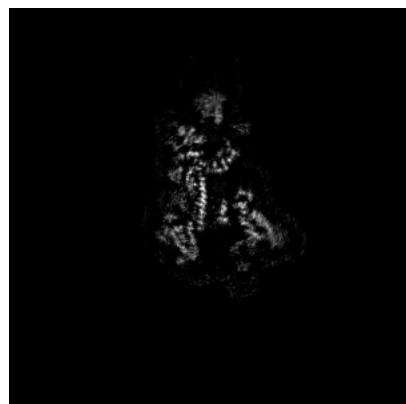


Z

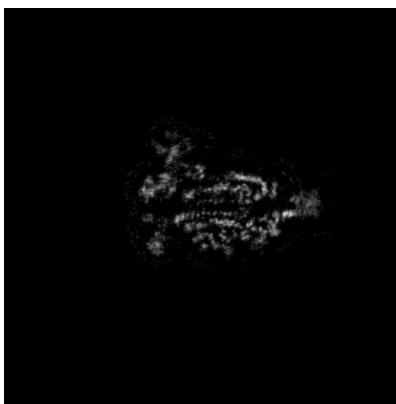
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 240



Y Index: 240

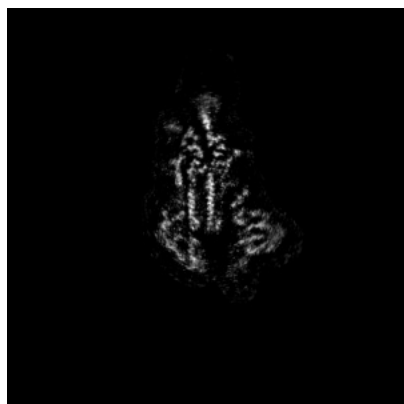


Z Index: 240

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

6.3 Largest variance slices [i](#)

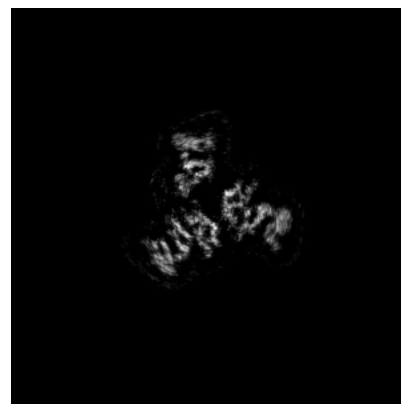
6.3.1 Primary map



X Index: 233



Y Index: 232

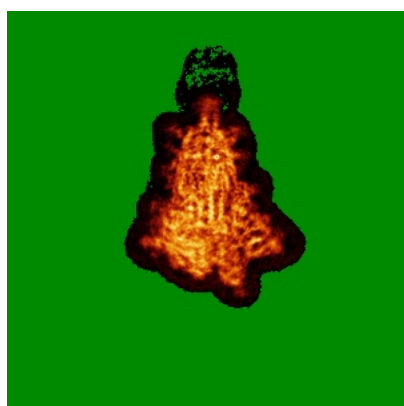


Z Index: 200

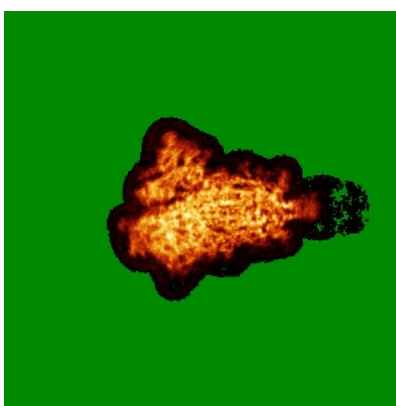
The images above show the largest variance slices of the map in three orthogonal directions.

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

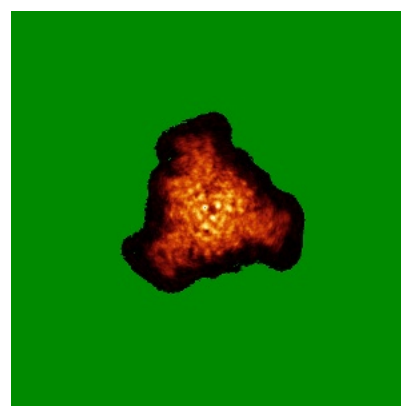
6.4.1 Primary map



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

This section was not generated.

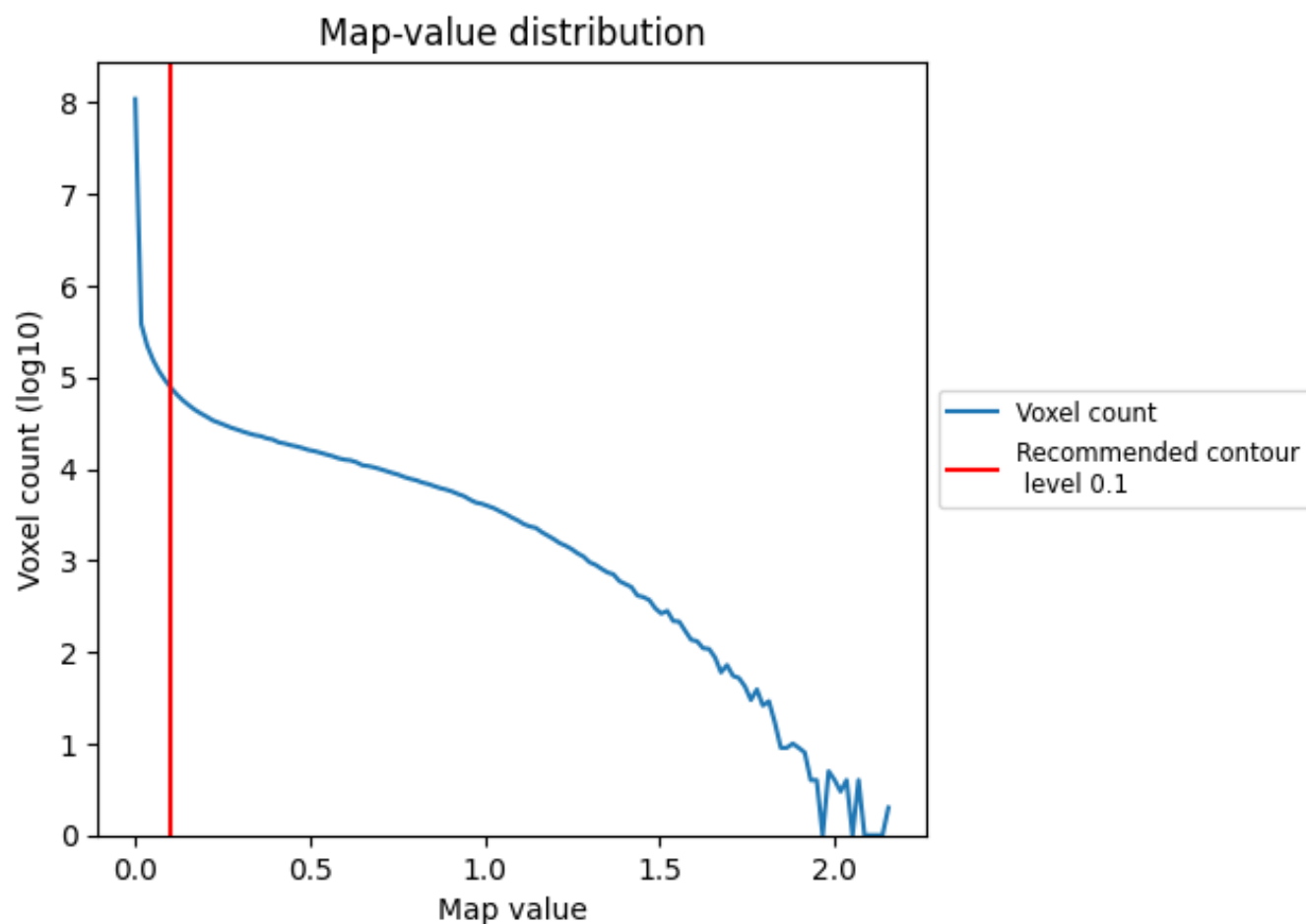
6.6 Mask visualisation

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

7 Map analysis [i](#)

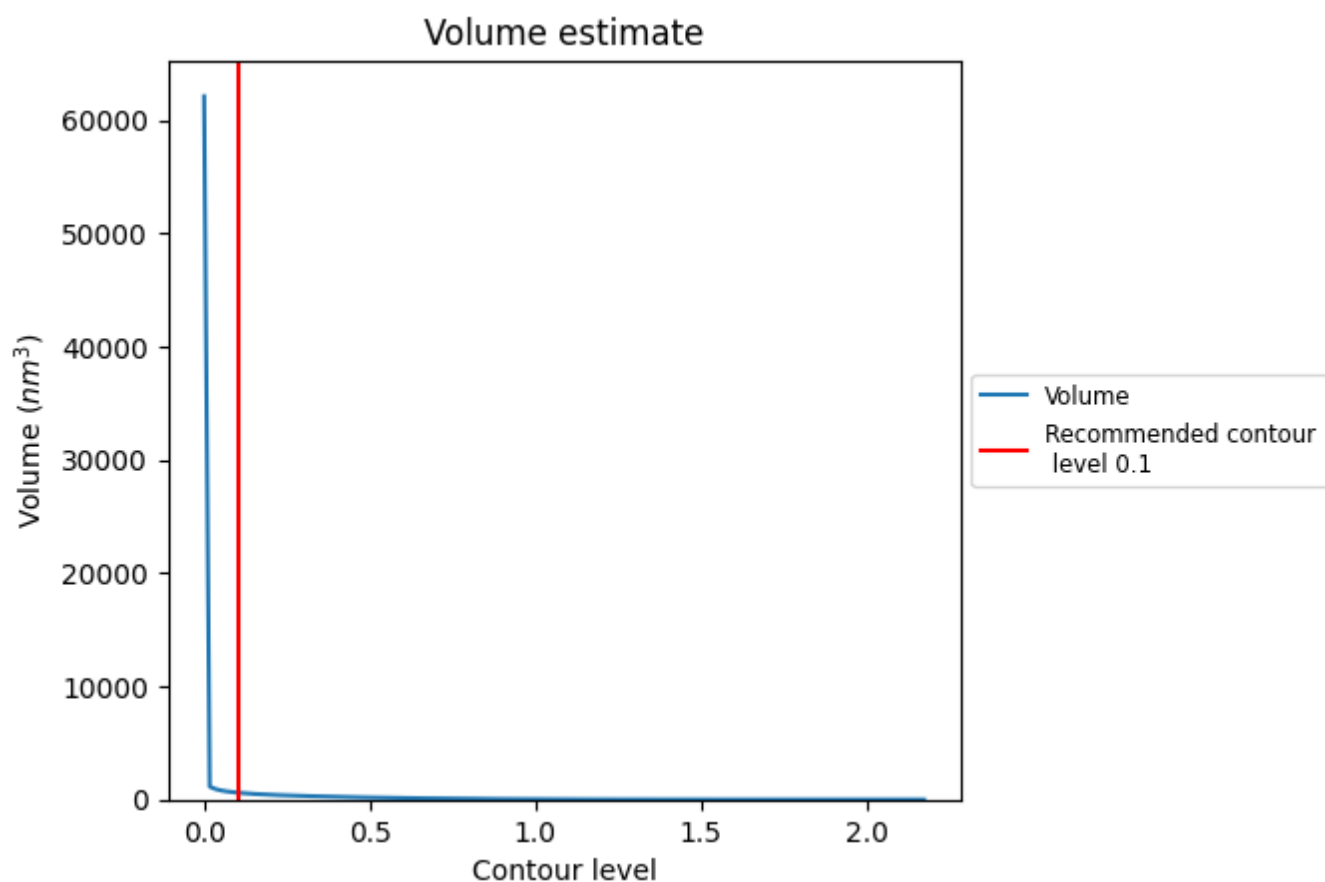
This section contains the results of statistical analysis of the map.

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.

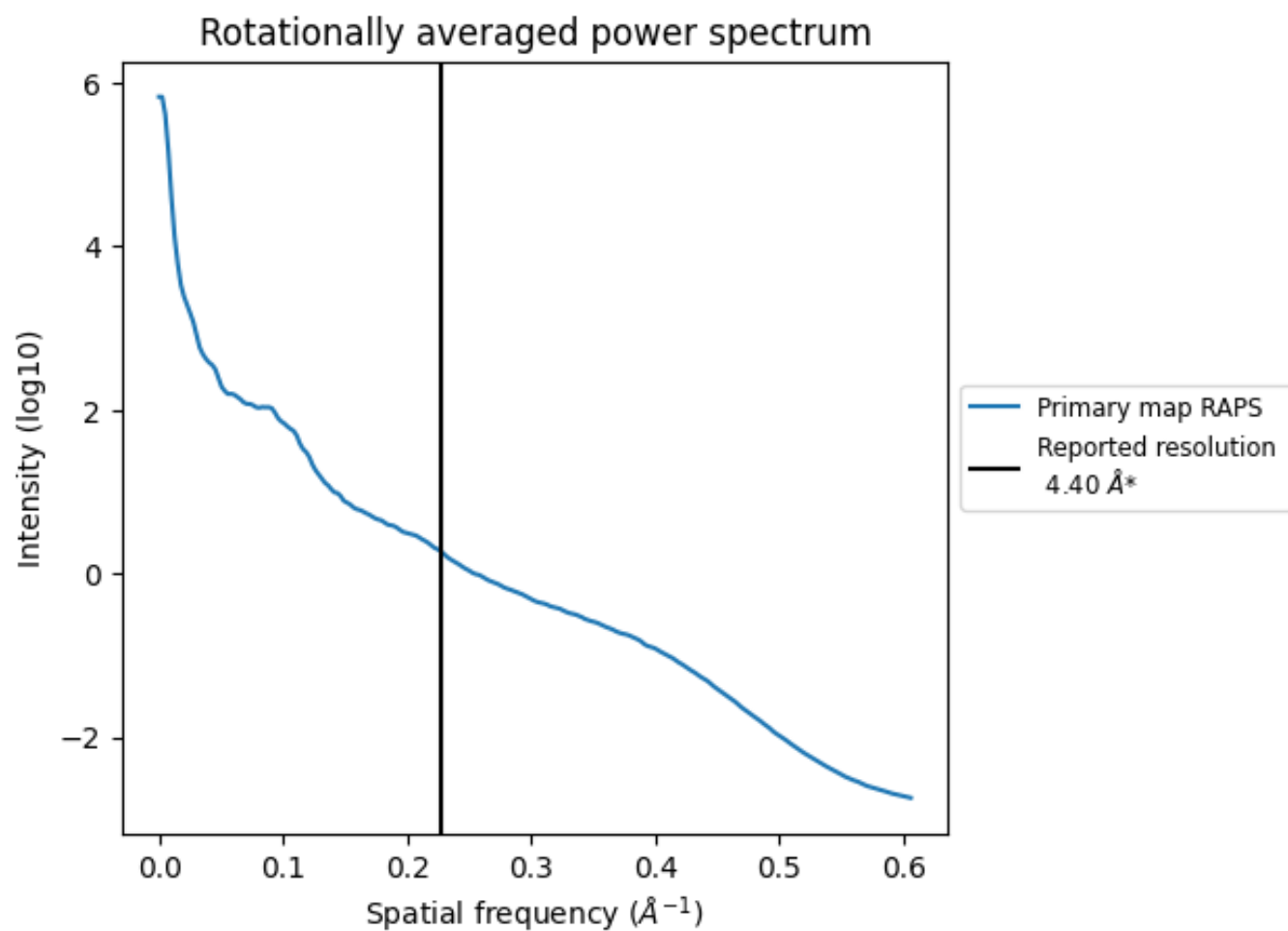
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 606 nm³; this corresponds to an approximate mass of 547 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 0.227 Å⁻¹

8 Fourier-Shell correlation

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

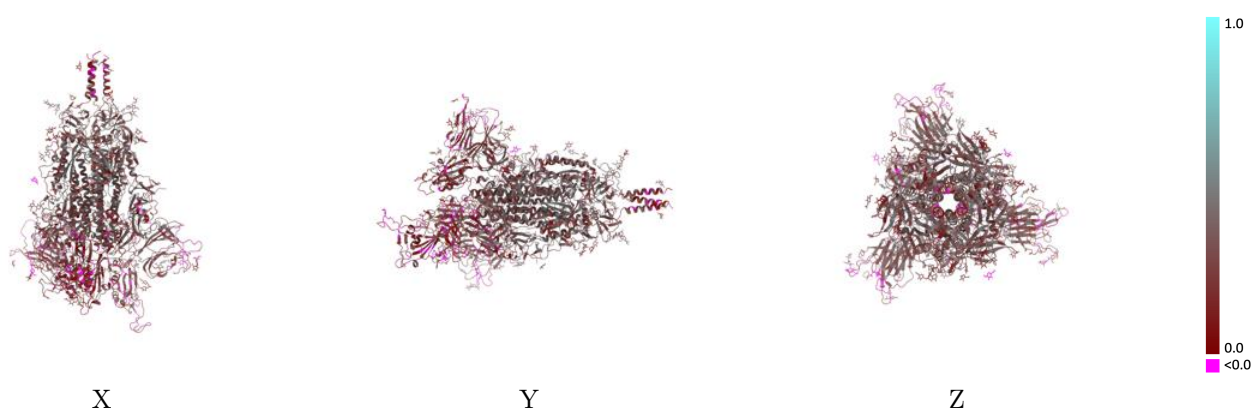
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-24988 and PDB model 7SBT. Per-residue inclusion information can be found in section 3 on page 13.

9.1 Map-model overlay [i](#)

This section was not generated.

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

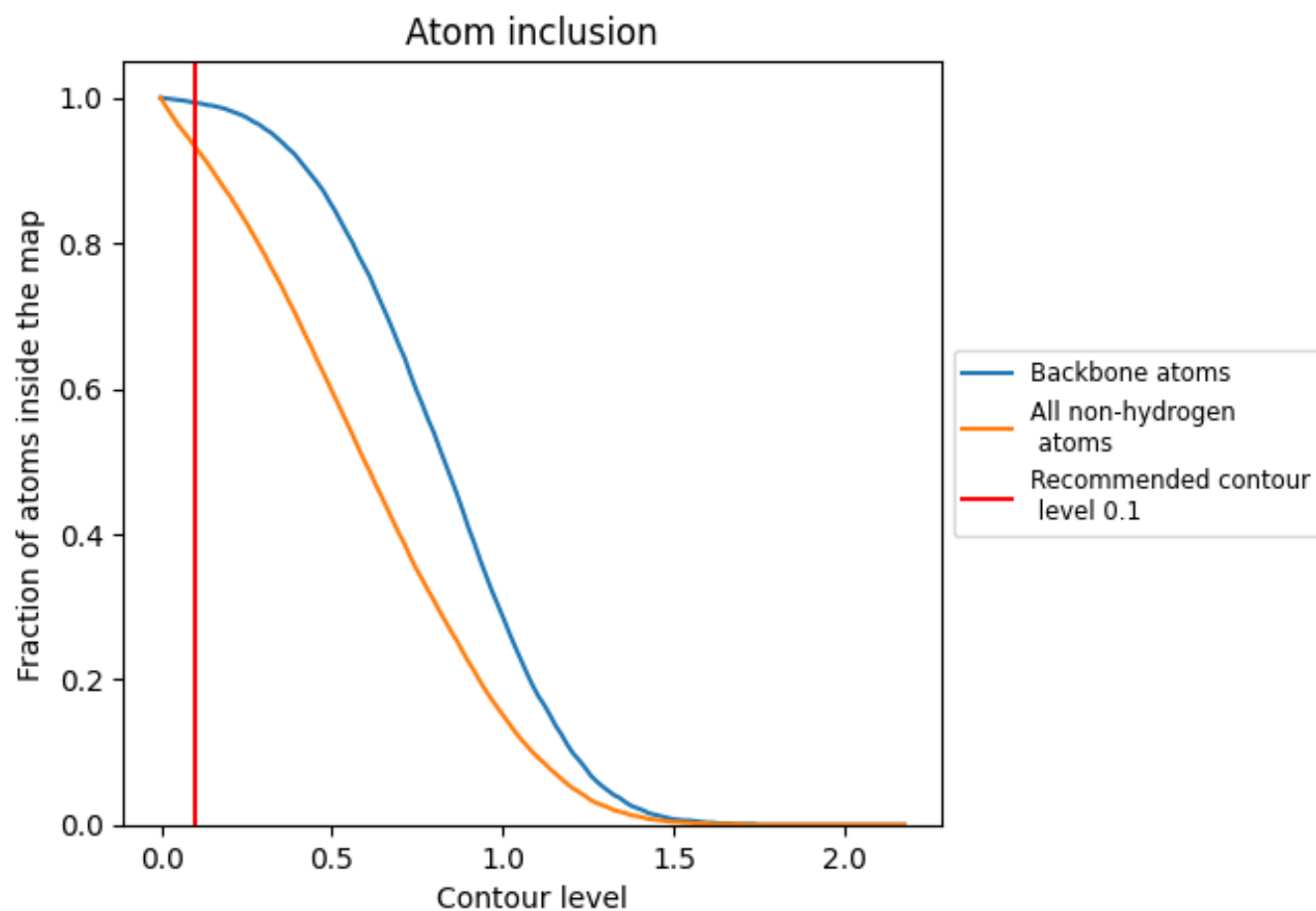


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 99% of all backbone atoms, 93% 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	 0.9320	 0.2690
A	 0.9270	 0.2670
B	 0.9370	 0.2750
C	 0.9330	 0.2690
D	 0.9290	 0.1660
E	 1.0000	 0.3870
F	 0.9290	 0.3430
G	 0.9290	 0.2400
H	 0.8950	 0.1340
I	 1.0000	 0.2030
J	 1.0000	 0.4100
K	 0.6920	 0.1390
L	 0.9290	 0.3000
M	 1.0000	 0.2140
N	 0.9640	 0.1860
O	 1.0000	 0.0860
P	 0.9640	 0.3480
Q	 0.9640	 0.2670
R	 0.8950	 0.1790
S	 0.9740	 0.2900
T	 1.0000	 0.2660
U	 0.8210	 0.0670
V	 0.9640	 0.0990
W	 1.0000	 0.2030
X	 0.9640	 0.2360
Y	 1.0000	 0.2160
Z	 0.9290	 0.2510
a	 0.8570	 0.2680
b	 1.0000	 0.3310
c	 0.8420	 0.1590
d	 1.0000	 0.2870
e	 1.0000	 0.3340
f	 0.7690	 0.0780

