



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 03:52 PM UTC

PDB ID : 8GJM / pdb_00008gjm
EMDB ID : EMD-40094
Title : 17b10 fab in complex with full-length SARS-CoV-2 Spike G614 trimer
Authors : Kwon, H.J.; Zhang, J.; Kosikova, M.; Tang, W.C.; Rodriguez, U.O.; Peng, H.Q.; Meseda, C.A.; Pedro, C.L.; Schmeisser, F.; Lu, J.M.; Zhou, B.; Davis, C.T.; Wentworth, D.E.; Chen, W.H.; Shriver, M.C.; Pasetti, M.F.; Weir, J.P.; Chen, B.; Xie, H.
Deposited on : 2023-03-16
Resolution : 2.80 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)

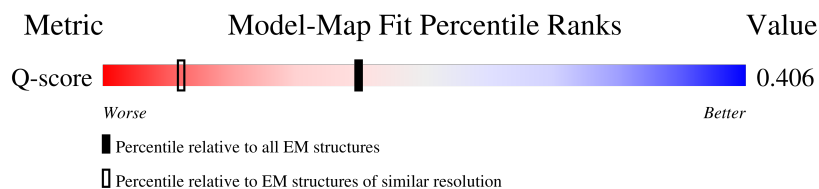
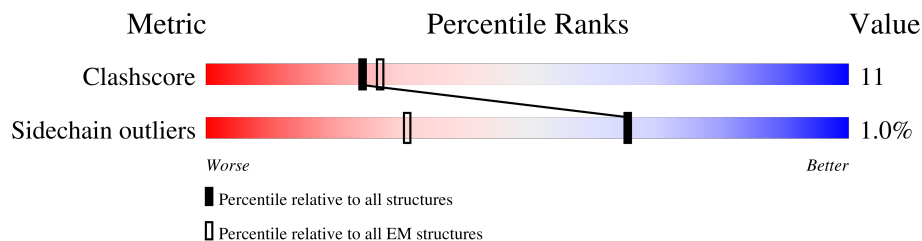
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11806 (2.30 - 3.30)

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	138	

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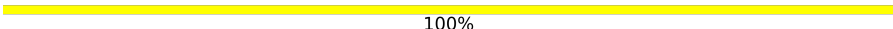
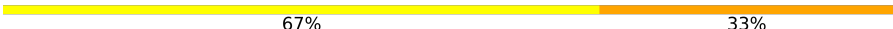
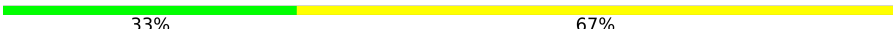
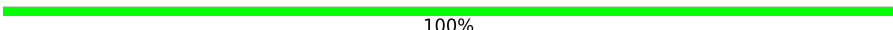
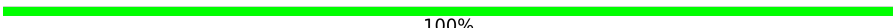
Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
3	E	127	
4	F	2	
4	G	2	
4	H	2	
4	I	2	
4	J	2	
4	K	2	
4	O	2	
4	P	2	
4	Q	2	
4	R	2	
4	S	2	
4	T	2	
4	U	2	
4	Y	2	
4	Z	2	
4	a	2	
4	b	2	
4	c	2	
4	d	2	
4	e	2	
4	f	2	
4	i	2	
4	j	2	
4	k	2	

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Mol	Chain	Length	Quality of chain
5	L	3	 67%33%
5	V	3	 100%
5	g	3	 67%33%
6	M	3	 33%67%
6	N	3	 100%
6	W	3	 67%33%
6	X	3	 100%
6	h	3	 33%67%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 28728 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	1089	Total	C	N	O	S	0	0
			8519	5438	1419	1623	39		
1	B	1088	Total	C	N	O	S	0	0
			8520	5441	1419	1622	38		
1	C	1099	Total	C	N	O	S	0	0
			8605	5493	1436	1637	39		

There are 114 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	614	GLY	ASP	conflict	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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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	614	GLY	ASP	conflict	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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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	614	GLY	ASP	conflict	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

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Chain	Residue	Modelled	Actual	Comment	Reference
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 a protein called Heavy chain of 17B10 fab.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	119	Total	C	N	O	S	0	0
			934	583	156	191	4		

- Molecule 3 is a protein called Light chain of 17B10 fab.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	106	Total	C	N	O	S	0	0
			819	516	134	166	3		

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



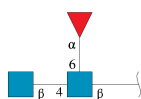
Mol	Chain	Residues	Atoms					AltConf	Trace
4	F	2	Total	C	N	O		0	0
			28	16	2	10			
4	G	2	Total	C	N	O		0	0
			28	16	2	10			
4	H	2	Total	C	N	O		0	0
			28	16	2	10			
4	I	2	Total	C	N	O		0	0
			28	16	2	10			
4	J	2	Total	C	N	O		0	0
			28	16	2	10			
4	K	2	Total	C	N	O		0	0
			28	16	2	10			
4	O	2	Total	C	N	O		0	0
			28	16	2	10			

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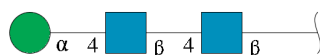
Mol	Chain	Residues	Atoms				AltConf	Trace
4	P	2	Total	C	N	O	0	0
			28	16	2	10		
4	Q	2	Total	C	N	O	0	0
			28	16	2	10		
4	R	2	Total	C	N	O	0	0
			28	16	2	10		
4	S	2	Total	C	N	O	0	0
			28	16	2	10		
4	T	2	Total	C	N	O	0	0
			28	16	2	10		
4	U	2	Total	C	N	O	0	0
			28	16	2	10		
4	Y	2	Total	C	N	O	0	0
			28	16	2	10		
4	Z	2	Total	C	N	O	0	0
			28	16	2	10		
4	a	2	Total	C	N	O	0	0
			28	16	2	10		
4	b	2	Total	C	N	O	0	0
			28	16	2	10		
4	c	2	Total	C	N	O	0	0
			28	16	2	10		
4	d	2	Total	C	N	O	0	0
			28	16	2	10		
4	e	2	Total	C	N	O	0	0
			28	16	2	10		
4	f	2	Total	C	N	O	0	0
			28	16	2	10		
4	i	2	Total	C	N	O	0	0
			28	16	2	10		
4	j	2	Total	C	N	O	0	0
			28	16	2	10		
4	k	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



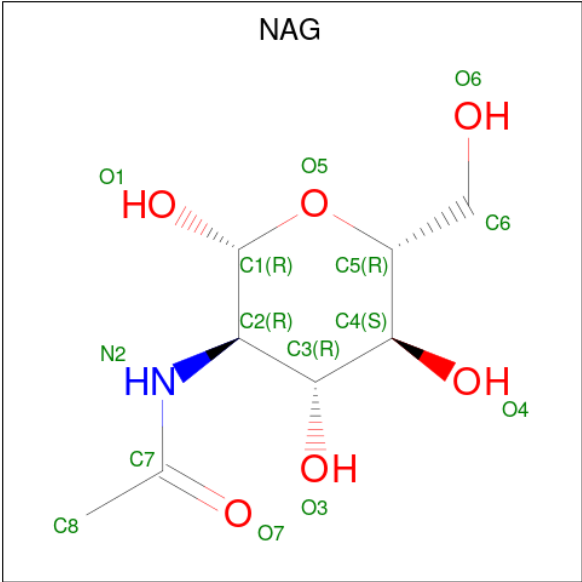
Mol	Chain	Residues	Atoms				AltConf	Trace
5	L	3	Total	C	N	O	0	0
			38	22	2	14		
5	V	3	Total	C	N	O	0	0
			38	22	2	14		
5	g	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 6 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
6	M	3	Total	C	N	O	0	0
			39	22	2	15		
6	N	3	Total	C	N	O	0	0
			39	22	2	15		
6	W	3	Total	C	N	O	0	0
			39	22	2	15		
6	X	3	Total	C	N	O	0	0
			39	22	2	15		
6	h	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



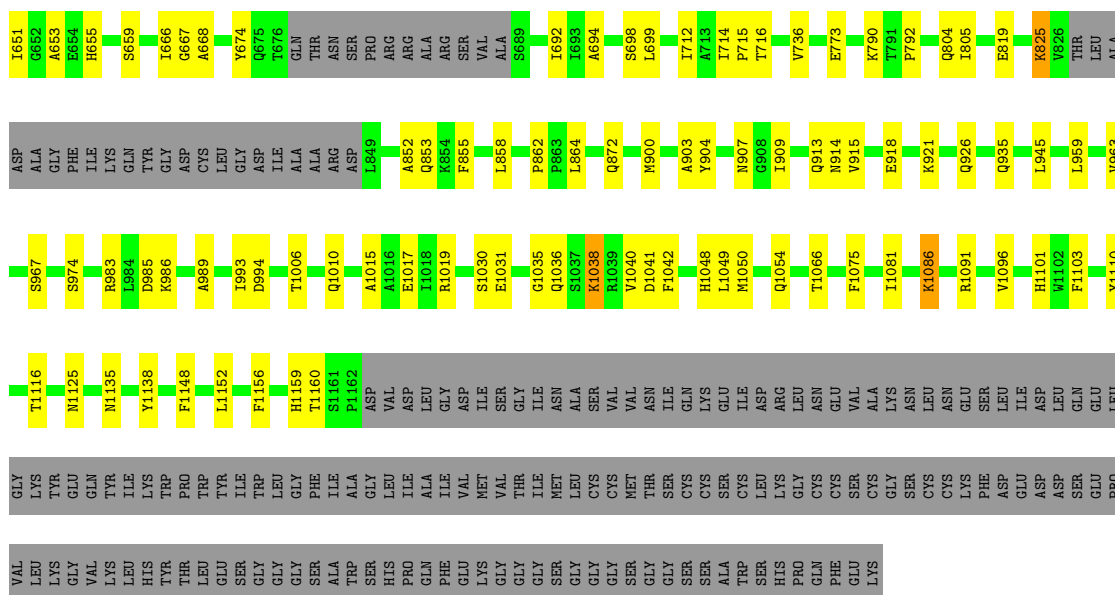
Mol	Chain	Residues	Atoms				AltConf
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	A	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7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

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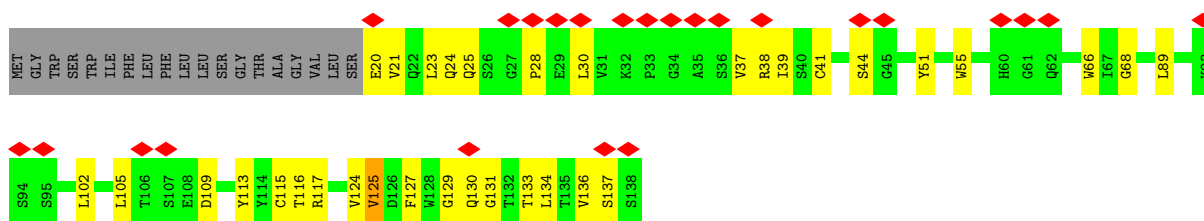
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Mol	Chain	Residues	Atoms				AltConf
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	C	1	Total	C	N	O	0
			14	8	1	5	
7	C	1	Total	C	N	O	0
			14	8	1	5	
7	C	1	Total	C	N	O	0
			14	8	1	5	
7	C	1	Total	C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





- Molecule 2: Heavy chain of 17B10 fab



- Molecule 3: Light chain of 17B10 fab



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



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

Chain G:  50% 50%

MAG1
MAG2

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

Chain H:  100%

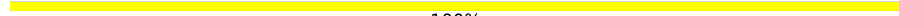
MAG1
MAG2

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

Chain I:  100%

MAG1
MAG2

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

Chain J:  100%

MAG1
MAG2

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

Chain K:  100%

MAG1
MAG2

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

Chain O:  50% 50%

MAG1
MAG2

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

Chain P:  50% 50%

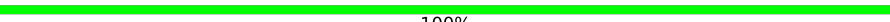
MAG1
MAG2

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

Chain Q:  50% 50%



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

Chain R:  100%



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

Chain S:  50% 50%

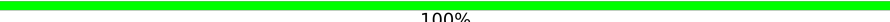


- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:  50% 50%



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

Chain U:  100%



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

Chain Y:  50% 50%



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

Chain Z:  100%

MAG1
MAG2

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

Chain a:  50% 50%

MAG1
MAG2

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

Chain b:  100%

MAG1
MAG2

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

Chain c:  100%

MAG1
MAG2

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

Chain d:  100%

MAG1
MAG2

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

Chain e:  100%

MAG1
MAG2

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

Chain f:  100%

MAG1
MAG2

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

Chain i:  100%

MAG1
MAG2

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

Chain j:  50% 50%

MAG1
MAG2

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

Chain k:  50% 50%

MAG1
MAG2

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

Chain L:  67% 33%

MAG1
MAG2
FUC3

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

Chain V:  100%

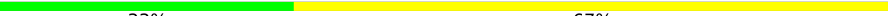
MAG1
MAG2
FUC3

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

Chain g:  67% 33%

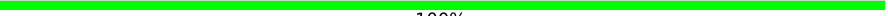
MAG1
MAG2
FUC3

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

Chain M:  33% 67%

MAG1
MAG2
MAN3

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

Chain N:  100%

MAG1
MAG2
MAN3

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

Chain W:  67% 33%

MAG1
MAG2
MAN3

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

Chain h:  33% 67%

MAG1
MAG2
MAN3

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	439265	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 ($e^-/\text{\AA}^2$)	54.442	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.520	Depositor
Minimum map value	-0.164	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.011	Depositor
Recommended contour level	0.02	Depositor
Map size (Å)	396.0, 396.0, 396.0	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.825, 0.825, 0.825	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.77	0/8715	1.18	6/11854 (0.1%)
1	B	0.79	0/8717	1.20	11/11858 (0.1%)
1	C	0.81	0/8806	1.20	4/11982 (0.0%)
2	D	0.78	0/954	1.15	0/1290
3	E	0.78	0/837	1.18	1/1130 (0.1%)
All	All	0.79	0/28029	1.19	22/38114 (0.1%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1091	ARG	N-CA-C	-7.40	102.81	111.03
1	B	103	GLY	N-CA-C	7.11	119.39	111.85
1	A	810	SER	N-CA-C	6.72	121.19	112.92
1	B	213	VAL	N-CA-C	6.28	116.32	110.42
3	E	81	ARG	N-CA-C	6.14	120.24	112.87
1	C	96	GLU	N-CA-C	6.02	118.00	108.79
1	A	1110	TYR	CA-C-N	5.94	128.42	122.00
1	A	1110	TYR	C-N-CA	5.94	128.42	122.00
1	B	520	ALA	N-CA-CB	-5.79	105.32	111.29
1	C	346	ARG	N-CA-C	5.77	118.11	108.02
1	C	95	THR	N-CA-C	-5.73	100.33	109.96
1	B	620	VAL	CA-C-N	5.50	125.02	119.19
1	B	620	VAL	C-N-CA	5.50	125.02	119.19
1	B	983	ARG	N-CA-C	5.33	118.94	112.23
1	A	524	VAL	CA-C-O	-5.27	118.05	122.63
1	B	1110	TYR	CA-C-N	5.16	126.82	122.28
1	B	1110	TYR	C-N-CA	5.16	126.82	122.28
1	A	619	GLU	CA-C-N	5.13	123.47	120.24
1	A	619	GLU	C-N-CA	5.13	123.47	120.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99	ASN	CA-C-N	5.12	127.11	120.56
1	B	99	ASN	C-N-CA	5.12	127.11	120.56
1	B	468	ILE	N-CA-C	-5.02	108.23	113.10

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	8519	0	8298	192	0
1	B	8520	0	8301	205	0
1	C	8605	0	8383	247	0
2	D	934	0	888	37	0
3	E	819	0	795	16	0
4	F	28	0	25	1	0
4	G	28	0	25	2	0
4	H	28	0	25	0	0
4	I	28	0	25	0	0
4	J	28	0	25	1	0
4	K	28	0	25	0	0
4	O	28	0	25	1	0
4	P	28	0	25	1	0
4	Q	28	0	25	4	0
4	R	28	0	25	0	0
4	S	28	0	25	1	0
4	T	28	0	25	0	0
4	U	28	0	25	0	0
4	Y	28	0	25	3	0
4	Z	28	0	25	0	0
4	a	28	0	25	4	0
4	b	28	0	25	2	0
4	c	28	0	25	0	0
4	d	28	0	25	0	0
4	e	28	0	25	1	0
4	f	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	i	28	0	25	0	0
4	j	28	0	25	1	0
4	k	28	0	25	1	0
5	L	38	0	34	0	0
5	V	38	0	34	1	0
5	g	38	0	34	1	0
6	M	39	0	34	2	0
6	N	39	0	34	0	0
6	W	39	0	34	1	0
6	X	39	0	34	0	0
6	h	39	0	34	3	0
7	A	126	0	117	1	0
7	B	112	0	104	0	0
7	C	112	0	104	2	0
All	All	28728	0	27862	602	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 11.

All (602) 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:B:102:ARG:NH1	1:B:179:LEU:HD23	1.25	1.50
1:B:102:ARG:HH12	1:B:179:LEU:CD2	1.25	1.45
1:B:102:ARG:NH1	1:B:179:LEU:CD2	1.84	1.34
1:B:102:ARG:CZ	1:B:179:LEU:CD2	2.25	1.14
1:B:102:ARG:NH2	1:B:179:LEU:CD2	2.14	1.10
1:B:102:ARG:HH22	1:B:179:LEU:CD2	1.67	1.06
1:B:14:GLN:OE1	1:B:158:ARG:NH1	1.87	1.06
2:D:102:LEU:HB3	2:D:105:LEU:HD21	1.39	1.04
1:A:457:ARG:NH2	1:A:467:ASP:OD2	1.96	0.96
1:B:102:ARG:NH2	1:B:179:LEU:HD21	1.78	0.95
1:B:102:ARG:HH22	1:B:179:LEU:HD22	1.31	0.95
1:B:102:ARG:HH12	1:B:179:LEU:HD22	1.30	0.94
1:A:403:ARG:HE	1:A:495:TYR:HD1	1.03	0.94
1:B:102:ARG:NH2	1:B:179:LEU:HD22	1.85	0.89
1:B:102:ARG:NH1	1:B:179:LEU:HD22	1.84	0.87
1:B:804:GLN:HE21	1:B:931:ILE:HG22	1.38	0.87
1:A:352:ALA:HB1	1:A:466:ARG:HH21	1.41	0.85
1:C:634:ARG:HG3	1:C:634:ARG:HH11	1.41	0.85
1:A:403:ARG:HG3	1:A:495:TYR:CE1	2.12	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:983:ARG:HD3	1:C:517:LEU:HD13	1.57	0.84
1:C:568:ASP:HB2	1:C:574:ASP:HB2	1.60	0.83
1:B:14:GLN:OE1	1:B:158:ARG:CZ	2.27	0.82
1:A:466:ARG:NE	1:A:468:ILE:HD11	1.96	0.80
1:A:1156:PHE:CE2	1:B:1155:TYR:HB3	2.16	0.80
1:B:589:PRO:HG3	1:C:855:PHE:HA	1.63	0.80
1:A:34:ARG:NH1	1:A:191:GLU:OE2	2.14	0.80
1:B:620:VAL:O	1:B:634:ARG:NH2	2.15	0.79
1:B:101:ILE:HB	1:B:190:ARG:HH22	1.47	0.79
1:A:720:ILE:HG12	1:A:926:GLN:NE2	1.97	0.79
1:A:352:ALA:HA	1:A:466:ARG:HE	1.47	0.79
3:E:24:MET:HB3	3:E:44:ARG:O	1.83	0.79
1:C:15:CYS:SG	1:C:136:CYS:CB	2.72	0.78
1:B:102:ARG:HH12	1:B:179:LEU:HD23	0.80	0.78
2:D:28:PRO:HA	2:D:133:THR:HB	1.66	0.78
1:C:326:ILE:HD11	1:C:534:VAL:HG22	1.63	0.78
1:B:102:ARG:HH22	1:B:179:LEU:HD21	1.42	0.77
1:A:200:TYR:OH	1:C:394:ASN:CB	2.34	0.76
1:B:390:LEU:HD21	1:C:983:ARG:HG2	1.68	0.76
1:A:395:VAL:HG22	1:A:515:PHE:CD1	2.21	0.76
2:D:30:LEU:HD11	2:D:137:SER:O	1.85	0.75
1:B:617:CYS:N	1:B:644:GLN:OE1	2.20	0.75
1:C:620:VAL:HG11	1:C:651:ILE:CD1	2.16	0.75
1:B:707:TYR:HB3	1:C:792:PRO:HG3	1.66	0.74
1:C:327:VAL:HG11	1:C:528:LYS:HG2	1.70	0.74
1:B:623:ALA:HB3	1:B:634:ARG:NH2	2.04	0.73
1:C:32:PHE:O	1:C:34:ARG:HG2	1.89	0.73
1:A:922:LEU:O	1:A:926:GLN:HG3	1.89	0.73
1:A:359:SER:HA	1:A:524:VAL:HG22	1.71	0.73
1:A:357:ARG:HH12	1:B:167:THR:HG23	1.53	0.72
1:C:110:LEU:HB3	1:C:135:PHE:CD2	2.23	0.72
1:A:403:ARG:NE	1:A:495:TYR:HD1	1.84	0.72
1:B:123:ALA:HB3	4:Q:1:NAG:H82	1.72	0.72
1:B:608:VAL:C	1:B:692:ILE:HD12	2.14	0.71
1:B:623:ALA:O	1:B:634:ARG:NE	2.23	0.71
1:C:21:ARG:HD3	1:C:79:PHE:HB3	1.72	0.71
1:C:559:PHE:HZ	1:C:566:GLY:HA3	1.55	0.71
1:A:403:ARG:NE	1:A:495:TYR:CD1	2.58	0.70
1:A:395:VAL:HG22	1:A:515:PHE:HD1	1.56	0.70
1:A:547:THR:OG1	1:B:978:ASN:ND2	2.25	0.70
1:B:468:ILE:HD11	4:k:1:NAG:H82	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:273:ARG:NH1	1:C:290:ASP:OD1	2.25	0.70
1:A:905:ARG:HH11	1:A:1036:GLN:HB2	1.57	0.69
1:C:903:ALA:HB1	1:C:913:GLN:HG3	1.72	0.69
1:A:852:ALA:HB1	1:C:570:ALA:HB2	1.74	0.69
1:B:89:GLY:HA3	1:B:270:LEU:HD12	1.72	0.69
1:A:466:ARG:HE	1:A:468:ILE:HD11	1.56	0.69
1:C:15:CYS:SG	1:C:136:CYS:HB3	2.31	0.69
1:A:1035:GLY:HA3	1:C:1040:VAL:HG21	1.75	0.69
1:A:530:SER:CB	1:A:580:GLN:HE22	2.06	0.69
2:D:124:VAL:HG12	2:D:125:VAL:HG23	1.74	0.68
1:C:565:PHE:HD1	1:C:576:VAL:HG23	1.57	0.68
1:A:230:PRO:HG3	1:C:357:ARG:NH1	2.09	0.67
1:B:312:ILE:HB	1:B:664:ILE:CG2	2.25	0.67
1:C:124:THR:HB	4:b:1:NAG:HN2	1.60	0.67
1:A:653:ALA:HA	1:A:692:ILE:O	1.95	0.67
1:B:562:PHE:HB2	1:C:41:LYS:HD2	1.77	0.66
1:C:611:LEU:HD22	1:C:666:ILE:HG23	1.77	0.66
3:E:52:TYR:HB2	3:E:112:ALA:HB2	1.76	0.66
1:C:89:GLY:HA3	1:C:270:LEU:HD12	1.78	0.66
3:E:27:SER:HB2	3:E:28:PRO:HD3	1.77	0.66
1:A:52:GLN:HA	1:A:274:THR:HA	1.78	0.66
1:A:530:SER:OG	1:A:580:GLN:NE2	2.29	0.66
1:B:149:ASN:HD22	1:B:153:MET:HE3	1.61	0.66
1:A:720:ILE:HG12	1:A:926:GLN:HE21	1.61	0.66
1:A:996:LEU:O	1:A:1000:ARG:HG3	1.96	0.66
1:C:630:THR:HB	1:C:631:PRO:HD3	1.76	0.65
1:B:600:PRO:HG3	1:B:674:TYR:CG	2.31	0.65
1:C:15:CYS:SG	4:a:1:NAG:H82	2.36	0.65
1:C:149:ASN:HD22	1:C:153:MET:HE3	1.61	0.65
1:B:339:GLY:HA2	4:Y:1:NAG:H81	1.78	0.65
1:B:1041:ASP:HB2	1:C:1030:SER:HB3	1.78	0.65
1:B:64:TRP:HE1	1:B:264:ALA:HB1	1.60	0.65
1:C:319:ARG:HB2	1:C:319:ARG:NH1	2.12	0.65
1:C:81:ASN:HB3	1:C:242:LEU:HD13	1.78	0.64
1:C:279:TYR:CE2	1:C:285:ILE:HG13	2.32	0.64
1:B:567:ARG:NE	1:C:42:VAL:HG21	2.13	0.64
1:A:125:ASN:HD21	4:G:1:NAG:H2	1.61	0.64
1:A:42:VAL:CG1	1:C:567:ARG:HB2	2.28	0.64
1:A:42:VAL:HG11	1:C:567:ARG:HB2	1.77	0.64
1:B:598:ILE:HD13	1:B:666:ILE:HD11	1.79	0.64
1:A:644:GLN:HA	1:A:644:GLN:HE21	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:600:PRO:HB3	1:C:674:TYR:CB	2.27	0.64
1:B:97:LYS:HD3	1:B:182:LYS:HB2	1.78	0.64
1:A:149:ASN:HD22	1:A:153:MET:HE3	1.61	0.64
1:A:96:GLU:O	1:A:188:ASN:ND2	2.31	0.64
1:B:146:HIS:CD2	1:B:150:LYS:HA	2.33	0.64
1:C:33:THR:HB	1:C:220:PHE:HD1	1.62	0.64
1:B:102:ARG:CZ	1:B:179:LEU:HD21	2.18	0.63
1:C:146:HIS:CD2	1:C:150:LYS:HA	2.33	0.63
1:C:147:LYS:HE3	7:C:1402:NAG:H81	1.80	0.63
1:A:147:LYS:HE3	7:A:1402:NAG:H81	1.80	0.63
1:A:862:PRO:HG3	1:C:647:ALA:HA	1.81	0.63
1:B:715:PRO:HA	1:B:1072:GLU:HA	1.79	0.63
1:C:819:GLU:HG3	1:C:1054:GLN:OE1	1.98	0.63
2:D:55:TRP:CD1	2:D:89:LEU:CD2	2.81	0.63
1:A:756:TYR:HB3	1:A:759:PHE:HD1	1.64	0.63
1:C:612:TYR:HB2	1:C:649:CYS:HB3	1.80	0.63
1:C:634:ARG:HH11	1:C:634:ARG:CG	2.11	0.63
2:D:28:PRO:CA	2:D:133:THR:HB	2.29	0.63
1:A:608:VAL:C	1:A:692:ILE:HD12	2.24	0.62
1:A:329:PHE:CD1	1:A:528:LYS:HB3	2.33	0.62
1:A:146:HIS:CD2	1:A:150:LYS:HA	2.33	0.62
1:A:905:ARG:NH1	1:A:1036:GLN:HB2	2.15	0.62
1:C:609:ALA:CB	1:C:653:ALA:HB2	2.29	0.62
1:B:102:ARG:CZ	1:B:179:LEU:HD22	2.10	0.62
1:A:57:PRO:HB3	1:A:273:ARG:NH2	2.15	0.62
1:C:137:ASN:HB3	4:a:1:NAG:C1	2.30	0.61
1:A:609:ALA:N	1:A:692:ILE:CD1	2.63	0.61
1:C:327:VAL:HA	1:C:542:ASN:HB3	1.82	0.61
1:A:352:ALA:HA	1:A:466:ARG:NE	2.15	0.61
1:B:620:VAL:HG11	1:B:651:ILE:CD1	2.29	0.61
1:A:646:ARG:HB2	1:A:668:ALA:CB	2.30	0.61
3:E:24:MET:SD	3:E:45:ALA:HA	2.41	0.61
1:A:983:ARG:HB3	1:C:382:VAL:HA	1.83	0.61
1:A:609:ALA:N	1:A:692:ILE:HD12	2.16	0.60
1:B:567:ARG:HG3	1:C:42:VAL:HG11	1.82	0.60
1:A:466:ARG:CZ	1:A:468:ILE:HD11	2.30	0.60
1:C:853:GLN:HG2	1:C:963:VAL:HG21	1.82	0.60
1:C:15:CYS:CB	1:C:136:CYS:HG	2.14	0.60
1:C:326:ILE:HB	1:C:541:PHE:HA	1.84	0.59
2:D:55:TRP:CD1	2:D:89:LEU:HD22	2.37	0.59
1:A:530:SER:CB	1:A:580:GLN:NE2	2.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:559:PHE:HB2	1:B:577:ARG:HH21	1.68	0.59
1:C:617:CYS:SG	1:C:644:GLN:HB2	2.41	0.59
3:E:37:GLU:N	3:E:37:GLU:OE1	2.36	0.59
1:C:634:ARG:HG2	1:C:634:ARG:O	2.01	0.59
1:A:127:VAL:HG11	4:G:1:NAG:H81	1.84	0.59
1:A:707:TYR:HB3	1:B:792:PRO:HG3	1.84	0.58
1:A:653:ALA:CB	1:A:692:ILE:HG22	2.33	0.58
1:A:983:ARG:HG2	1:C:390:LEU:HD21	1.85	0.58
1:A:409:GLN:HG2	1:A:414:GLN:OE1	2.03	0.58
1:A:352:ALA:HB1	1:A:466:ARG:NH2	2.16	0.58
1:A:455:LEU:HD22	1:A:493:GLN:HE21	1.68	0.58
2:D:55:TRP:CH2	2:D:115:CYS:HB3	2.38	0.58
3:E:81:ARG:NH2	3:E:99:GLU:HB2	2.19	0.58
1:B:1130:ILE:HD11	1:C:921:LYS:HE3	1.86	0.58
1:A:395:VAL:CG2	1:A:515:PHE:CD1	2.87	0.57
1:C:907:ASN:HD21	1:C:913:GLN:HB2	1.69	0.57
1:A:276:LEU:HB3	1:A:289:VAL:HB	1.86	0.57
1:A:409:GLN:HA	1:A:414:GLN:HG3	1.85	0.57
1:B:973:ILE:HD12	1:B:983:ARG:NH2	2.19	0.57
2:D:130:GLN:O	2:D:130:GLN:HG2	2.04	0.57
1:C:620:VAL:HG11	1:C:651:ILE:HD11	1.86	0.57
1:B:704:SER:HB2	1:C:790:LYS:HE3	1.85	0.57
1:A:357:ARG:NH1	1:B:167:THR:HG23	2.19	0.57
1:B:83:VAL:HG22	1:B:237:ARG:HG2	1.87	0.57
1:C:974:SER:HB2	1:C:983:ARG:HH22	1.70	0.57
1:B:647:ALA:HA	1:C:862:PRO:HG3	1.84	0.57
2:D:55:TRP:NE1	2:D:89:LEU:HD22	2.20	0.57
1:A:342:PHE:HB2	4:O:1:NAG:H82	1.87	0.57
1:A:654:GLU:N	1:A:692:ILE:O	2.31	0.57
1:A:667:GLY:HA2	1:B:864:LEU:HA	1.87	0.57
1:B:299:THR:CG2	1:B:597:VAL:HB	2.35	0.57
1:A:1017:GLU:HG2	1:B:1019:ARG:HH22	1.70	0.57
1:A:563:GLN:HE22	1:B:43:PHE:HD2	1.53	0.57
1:C:105:ILE:HD12	1:C:135:PHE:CE1	2.39	0.57
1:C:376:THR:HB	1:C:435:ALA:HB3	1.87	0.57
2:D:23:LEU:HB3	2:D:25:GLN:OE1	2.05	0.57
1:A:644:GLN:HA	1:A:644:GLN:NE2	2.20	0.56
1:C:559:PHE:CZ	1:C:566:GLY:HA3	2.38	0.56
1:A:29:THR:HB	1:A:62:VAL:HG23	1.86	0.56
1:A:983:ARG:CD	1:C:517:LEU:HD13	2.33	0.56
1:B:605:SER:HB2	1:B:674:TYR:CZ	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:273:ARG:NH1	1:C:290:ASP:OD2	2.38	0.56
1:A:1019:ARG:NH2	1:C:1017:GLU:HG2	2.21	0.56
3:E:76:SER:H	3:E:78:VAL:HG12	1.70	0.56
1:B:303:LEU:HD11	1:B:313:TYR:CD1	2.41	0.56
1:C:1101:HIS:CG	6:h:1:NAG:H5	2.41	0.56
1:C:635:VAL:HG22	1:C:635:VAL:O	2.05	0.56
1:B:605:SER:HB2	1:B:674:TYR:CE2	2.41	0.56
2:D:25:GLN:OE1	2:D:41:CYS:SG	2.64	0.56
1:B:15:CYS:O	4:P:1:NAG:H82	2.05	0.56
1:B:96:GLU:HA	1:B:186:PHE:CD1	2.41	0.55
1:B:118:LEU:CD1	1:B:159:VAL:HG11	2.36	0.55
1:B:1082:CYS:HB2	1:B:1132:ILE:HG12	1.87	0.55
1:C:49:HIS:HB3	1:C:279:TYR:HE1	1.72	0.55
1:A:609:ALA:CB	1:A:692:ILE:HG21	2.36	0.55
1:B:383:SER:HB2	1:C:985:ASP:H	1.71	0.55
1:C:612:TYR:O	1:C:649:CYS:N	2.40	0.55
1:A:200:TYR:OH	1:C:394:ASN:CG	2.50	0.55
1:A:498:GLN:H	1:A:501:ASN:ND2	2.05	0.55
1:A:1156:PHE:CE2	1:B:1155:TYR:CB	2.87	0.55
2:D:113:TYR:O	2:D:131:GLY:HA2	2.07	0.55
1:A:559:PHE:HB3	1:A:563:GLN:HB2	1.87	0.55
1:C:190:ARG:HD3	1:C:207:HIS:HD2	1.72	0.54
1:C:273:ARG:NH1	1:C:290:ASP:CG	2.64	0.54
2:D:23:LEU:HB2	2:D:129:GLY:HA2	1.90	0.54
1:A:487:ASN:HA	1:A:489:TYR:CE2	2.42	0.54
1:B:1106:GLN:HB2	1:B:1109:PHE:O	2.06	0.54
1:C:598:ILE:HD13	1:C:666:ILE:HD11	1.89	0.54
1:C:609:ALA:HB1	1:C:653:ALA:HB2	1.88	0.54
1:A:530:SER:HB3	1:A:580:GLN:HE22	1.72	0.54
1:B:123:ALA:HA	1:B:177:MET:CE	2.37	0.54
1:A:208:THR:HG22	1:A:210:ILE:HG23	1.90	0.54
1:A:332:ILE:HD13	1:A:527:PRO:HB2	1.89	0.54
1:A:479:PRO:HD3	3:E:52:TYR:CD2	2.43	0.54
2:D:21:VAL:HB	2:D:127:PHE:HD2	1.73	0.54
1:C:1031:GLU:HB2	1:C:1042:PHE:HE2	1.73	0.54
1:B:65:PHE:CZ	1:B:84:LEU:HD11	2.43	0.54
1:B:106:PHE:CD2	1:B:238:PHE:HB2	2.43	0.53
1:B:133:PHE:HB2	1:B:135:PHE:CE2	2.44	0.53
1:B:1031:GLU:HB2	1:B:1042:PHE:HE2	1.73	0.53
1:C:659:SER:HB3	1:C:698:SER:HB3	1.90	0.53
2:D:129:GLY:C	2:D:131:GLY:H	2.16	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:279:TYR:CE2	1:C:285:ILE:CG1	2.91	0.53
1:A:43:PHE:HB3	1:C:559:PHE:CZ	2.43	0.53
1:B:312:ILE:HB	1:B:664:ILE:HG21	1.90	0.53
1:B:972:ALA:HB2	1:B:995:ARG:HB3	1.89	0.53
1:B:611:LEU:HD22	1:B:666:ILE:HG23	1.91	0.53
2:D:28:PRO:HB3	2:D:133:THR:HB	1.89	0.53
1:B:714:ILE:HD12	1:B:1096:VAL:HG11	1.91	0.53
1:A:402:ILE:HD11	1:A:407:VAL:HA	1.91	0.53
1:B:949:GLN:O	1:B:953:ASN:OD1	2.26	0.53
1:C:189:LEU:HB2	1:C:210:ILE:HD13	1.91	0.53
1:A:1156:PHE:O	1:A:1160:THR:HG23	2.09	0.53
1:C:600:PRO:HB3	1:C:674:TYR:HB3	1.90	0.53
1:B:577:ARG:HE	1:B:584:ILE:HD11	1.75	0.52
1:A:521:PRO:HD3	1:B:199:GLY:HA3	1.91	0.52
1:C:599:THR:C	1:C:601:GLY:H	2.16	0.52
1:B:559:PHE:HB3	1:B:563:GLN:HB2	1.91	0.52
1:C:203:ILE:HG22	1:C:226:LEU:HD21	1.92	0.52
2:D:55:TRP:CD1	2:D:89:LEU:HD21	2.45	0.52
1:A:200:TYR:OH	1:C:394:ASN:HB3	2.07	0.52
1:B:716:THR:HG21	1:B:1073:LYS:HD3	1.91	0.52
1:A:102:ARG:HD3	1:A:121:ASN:O	2.09	0.52
1:B:32:PHE:HB2	1:B:218:GLN:HA	1.92	0.52
1:B:567:ARG:HD3	1:B:571:ASP:OD2	2.10	0.52
1:B:605:SER:CB	1:B:674:TYR:CZ	2.93	0.52
1:B:896:ILE:HD11	1:B:900:MET:HG2	1.91	0.52
1:B:396:TYR:CZ	1:C:230:PRO:HB3	2.45	0.52
1:C:575:ALA:HA	1:C:585:LEU:O	2.09	0.52
2:D:28:PRO:CB	2:D:133:THR:HB	2.40	0.52
1:A:478:THR:HG21	3:E:114:TYR:CE2	2.44	0.52
1:B:667:GLY:HA2	1:C:864:LEU:HA	1.92	0.52
1:A:567:ARG:HG3	1:B:42:VAL:CG1	2.39	0.52
1:B:557:LYS:HD3	1:C:43:PHE:CE2	2.44	0.52
1:C:1048:HIS:HA	1:C:1066:THR:HG22	1.91	0.52
1:A:653:ALA:HB2	1:A:692:ILE:HG22	1.92	0.51
1:B:1093:GLY:H	1:B:1107:ARG:HH11	1.57	0.51
1:C:615:VAL:HG12	1:C:649:CYS:HB2	1.92	0.51
1:C:328:ARG:N	1:C:542:ASN:O	2.38	0.51
1:C:909:ILE:HG22	1:C:1038:LYS:HE2	1.90	0.51
1:C:15:CYS:HG	1:C:136:CYS:CB	2.18	0.51
1:A:655:HIS:HA	1:A:694:ALA:O	2.11	0.51
1:A:872:GLN:HB2	1:C:699:LEU:CD1	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:804:GLN:HB3	1:C:935:GLN:OE1	2.10	0.51
1:C:853:GLN:HB3	1:C:858:LEU:HD12	1.92	0.51
1:A:560:LEU:HD12	1:A:562:PHE:HE1	1.74	0.51
1:C:565:PHE:CD1	1:C:576:VAL:HG23	2.42	0.51
1:B:608:VAL:HA	1:B:692:ILE:HD11	1.92	0.51
1:A:853:GLN:HB3	1:A:858:LEU:HB2	1.93	0.51
1:B:299:THR:HG21	1:B:597:VAL:HB	1.91	0.51
1:B:1107:ARG:HB3	1:C:904:TYR:CE2	2.45	0.51
1:C:858:LEU:HD13	1:C:959:LEU:HD22	1.93	0.51
1:B:310:LYS:HD3	1:B:663:ASP:OD1	2.11	0.51
1:C:634:ARG:HG2	1:C:637:SER:OG	2.11	0.51
1:B:737:ASP:HB2	1:B:740:MET:HB3	1.91	0.51
1:C:120:VAL:HG13	1:C:127:VAL:HG13	1.93	0.51
1:C:1006:THR:O	1:C:1010:GLN:HG2	2.11	0.51
1:C:674:TYR:HD1	1:C:692:ILE:HD13	1.76	0.50
1:C:611:LEU:CD2	1:C:666:ILE:HG23	2.42	0.50
1:A:560:LEU:H	1:A:563:GLN:HE21	1.59	0.50
1:C:634:ARG:CG	1:C:634:ARG:NH1	2.72	0.50
2:D:39:ILE:HD11	2:D:134:LEU:HD22	1.93	0.50
1:C:95:THR:HG22	1:C:186:PHE:CD2	2.46	0.50
1:B:609:ALA:N	1:B:692:ILE:HD12	2.26	0.50
1:B:342:PHE:HB2	4:Y:1:NAG:H82	1.94	0.50
1:B:569:ILE:HD12	1:B:569:ILE:H	1.77	0.50
1:A:1152:LEU:HD13	1:B:1152:LEU:HD23	1.94	0.50
1:C:129:LYS:HB3	1:C:133:PHE:HZ	1.76	0.50
1:C:196:ASN:HB2	1:C:201:PHE:HD1	1.77	0.50
1:A:540:ASN:HD22	1:A:549:THR:HG22	1.76	0.49
1:A:654:GLU:O	1:A:693:ILE:HA	2.12	0.49
1:B:105:ILE:HD12	1:B:135:PHE:CE1	2.47	0.49
1:A:1017:GLU:HG2	1:B:1019:ARG:NH2	2.27	0.49
1:C:314:GLN:HE21	1:C:613:GLN:HE22	1.58	0.49
1:C:1156:PHE:O	1:C:1160:THR:HG23	2.12	0.49
3:E:51:GLY:HA2	3:E:91:TYR:HE2	1.77	0.49
2:D:37:VAL:HG12	2:D:102:LEU:HB2	1.94	0.49
1:A:95:THR:O	1:A:186:PHE:CD1	2.66	0.49
1:C:559:PHE:CD1	1:C:584:ILE:HG21	2.48	0.49
1:A:709:ASN:HB3	1:B:796:ASP:HB3	1.94	0.49
1:C:102:ARG:HH22	1:C:179:LEU:HG	1.76	0.49
1:B:567:ARG:HG3	1:C:42:VAL:CG1	2.43	0.49
1:C:57:PRO:HG3	1:C:273:ARG:HD2	1.94	0.49
1:A:478:THR:HG21	3:E:114:TYR:CZ	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:M:1:NAG:H61	6:M:2:NAG:H82	1.94	0.49
1:A:926:GLN:HG2	4:J:2:NAG:H82	1.95	0.48
1:C:332:ILE:HB	1:C:362:VAL:HG23	1.95	0.48
1:B:97:LYS:C	1:B:98:SER:HG	2.21	0.48
1:C:926:GLN:HG2	4:e:2:NAG:H82	1.95	0.48
1:B:567:ARG:HE	1:C:42:VAL:HG21	1.77	0.48
1:B:1082:CYS:HB3	1:B:1134:ASN:OD1	2.12	0.48
1:A:567:ARG:HG3	1:B:42:VAL:HG13	1.96	0.48
1:A:1155:TYR:HB3	1:C:1156:PHE:CE2	2.49	0.48
2:D:51:TYR:CG	2:D:117:ARG:HD2	2.48	0.48
1:A:1019:ARG:HH22	1:C:1017:GLU:HG2	1.78	0.48
1:B:902:MET:HB3	1:B:916:LEU:HD11	1.96	0.48
5:g:2:NAG:H3	5:g:3:FUC:H61	1.95	0.48
1:A:457:ARG:HH22	1:A:467:ASP:CG	2.11	0.48
1:A:979:ASP:OD1	1:C:518:LEU:CD1	2.62	0.48
1:B:570:ALA:HB2	1:C:963:VAL:HG13	1.95	0.48
1:B:923:ILE:HA	1:B:926:GLN:HE21	1.78	0.48
1:B:1040:VAL:HG21	1:C:1035:GLY:HA3	1.94	0.48
1:A:608:VAL:C	1:A:692:ILE:CD1	2.87	0.48
1:B:609:ALA:N	1:B:692:ILE:CD1	2.77	0.48
1:A:395:VAL:CG2	1:A:515:PHE:HD1	2.25	0.48
1:B:608:VAL:C	1:B:692:ILE:CD1	2.86	0.48
1:B:973:ILE:HD12	1:B:983:ARG:HH21	1.78	0.48
1:B:16:VAL:HG21	1:B:158:ARG:HH11	1.79	0.47
1:C:56:LEU:HD22	1:C:91:TYR:CD2	2.49	0.47
1:C:714:ILE:HD12	1:C:1096:VAL:HG11	1.95	0.47
1:B:560:LEU:HB2	1:B:563:GLN:HG3	1.97	0.47
1:B:609:ALA:CB	1:B:653:ALA:HB2	2.44	0.47
1:A:147:LYS:HD2	1:A:147:LYS:HA	1.66	0.47
1:A:1155:TYR:HB3	1:C:1156:PHE:CZ	2.49	0.47
1:C:137:ASN:CG	4:a:1:NAG:H5	2.39	0.47
1:C:147:LYS:HD2	1:C:147:LYS:HA	1.66	0.47
1:C:1116:THR:HG22	1:C:1138:TYR:HB3	1.95	0.47
1:A:95:THR:O	1:A:186:PHE:CG	2.67	0.47
1:B:300:LYS:HE2	1:B:306:PHE:O	2.14	0.47
1:A:720:ILE:H	1:A:926:GLN:NE2	2.11	0.47
1:A:874:THR:HA	1:A:877:LEU:HD12	1.97	0.47
1:B:124:THR:HB	4:Q:1:NAG:HN2	1.80	0.47
1:B:141:LEU:O	1:B:243:ALA:HA	2.14	0.47
1:C:327:VAL:H	1:C:531:THR:HB	1.79	0.47
1:A:457:ARG:CD	1:A:459:SER:O	2.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:23:LEU:O	2:D:25:GLN:NE2	2.45	0.47
3:E:25:THR:O	3:E:44:ARG:HG2	2.14	0.47
1:A:313:TYR:O	1:A:596:SER:HA	2.15	0.47
1:C:600:PRO:HB2	1:C:604:THR:HB	1.97	0.47
1:C:1081:ILE:HD12	1:C:1135:ASN:HB3	1.97	0.47
1:A:565:PHE:HA	1:A:576:VAL:HA	1.96	0.46
1:C:321:GLN:HB2	1:C:628:GLN:O	2.15	0.46
1:A:659:SER:HB3	1:A:698:SER:HB3	1.96	0.46
1:A:1040:VAL:HG21	1:B:1035:GLY:HA3	1.97	0.46
1:B:289:VAL:HG23	1:B:306:PHE:CE2	2.51	0.46
1:C:68:ILE:O	1:C:69:HIS:C	2.58	0.46
1:B:615:VAL:H	1:B:648:GLY:HA2	1.81	0.46
1:C:989:ALA:O	1:C:993:ILE:HG12	2.14	0.46
1:A:756:TYR:HB3	1:A:759:PHE:CD1	2.48	0.46
1:C:625:HIS:O	1:C:626:ALA:C	2.59	0.46
2:D:102:LEU:HD13	2:D:134:LEU:HD21	1.97	0.46
1:B:620:VAL:HG11	1:B:651:ILE:HD11	1.97	0.46
1:B:634:ARG:HA	1:B:634:ARG:HD2	1.75	0.46
1:C:213:VAL:HG13	1:C:214:ARG:HG2	1.98	0.46
1:B:97:LYS:HE3	1:B:97:LYS:HB3	1.80	0.46
1:C:28:TYR:CE2	7:C:1401:NAG:H5	2.50	0.46
1:C:121:ASN:ND2	1:C:126:VAL:HG22	2.30	0.46
1:A:230:PRO:HG3	1:C:357:ARG:HH11	1.77	0.46
1:C:610:VAL:O	1:C:651:ILE:N	2.33	0.46
1:A:466:ARG:HE	1:A:468:ILE:CD1	2.28	0.46
1:A:466:ARG:NH2	1:A:468:ILE:HD11	2.30	0.46
1:B:987:VAL:HG13	1:C:413:GLY:HA3	1.98	0.46
1:B:1161:SER:HB2	1:B:1162:PRO:HD3	1.98	0.46
2:D:55:TRP:CZ3	2:D:115:CYS:HB3	2.51	0.46
1:B:1101:HIS:CG	6:W:1:NAG:H5	2.51	0.46
1:C:773:GLU:OE2	1:C:1015:ALA:HB1	2.15	0.46
1:A:206:LYS:HB3	1:A:223:LEU:HD22	1.98	0.45
1:B:147:LYS:HD2	1:B:147:LYS:HA	1.66	0.45
1:C:28:TYR:HB3	1:C:61:ASN:OD1	2.16	0.45
1:C:137:ASN:CB	4:a:1:NAG:C1	2.95	0.45
1:B:1077:THR:CG2	1:C:900:MET:HE3	2.46	0.45
1:B:1159:HIS:HB3	1:C:1159:HIS:CE1	2.51	0.45
1:C:319:ARG:HB2	1:C:319:ARG:HH11	1.79	0.45
1:A:601:GLY:C	1:A:603:ASN:N	2.72	0.45
1:B:392:PHE:CE2	1:B:517:LEU:HD21	2.51	0.45
1:B:624:ILE:HA	1:B:637:SER:OG	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:655:HIS:HA	1:B:694:ALA:O	2.16	0.45
1:B:970:PHE:O	1:B:995:ARG:HG2	2.16	0.45
1:C:634:ARG:HG3	1:C:634:ARG:NH1	2.20	0.45
1:C:712:ILE:HG22	1:C:1075:PHE:O	2.16	0.45
1:B:330:PRO:HD2	1:B:525:CYS:SG	2.57	0.45
1:B:1106:GLN:NE2	1:B:1111:GLU:OE2	2.49	0.45
1:C:327:VAL:O	1:C:531:THR:N	2.48	0.45
1:A:985:ASP:O	1:A:986:LYS:C	2.60	0.45
1:A:1159:HIS:C	1:A:1161:SER:H	2.25	0.45
1:B:736:VAL:HG13	1:B:858:LEU:HD23	1.98	0.45
1:C:34:ARG:O	1:C:56:LEU:HD23	2.15	0.45
1:A:365:TYR:HH	1:A:515:PHE:HE1	1.63	0.45
1:C:19:THR:O	1:C:21:ARG:N	2.49	0.45
1:A:179:LEU:H	1:A:179:LEU:HG	1.55	0.45
1:B:699:LEU:HD12	1:C:872:GLN:CB	2.47	0.45
1:B:1007:TYR:O	1:B:1011:GLN:HG2	2.17	0.45
1:C:320:VAL:HA	1:C:629:LEU:HG	1.99	0.45
1:C:610:VAL:O	1:C:650:LEU:HD12	2.16	0.45
1:C:825:LYS:HB2	1:C:945:LEU:HD22	1.99	0.45
1:A:490:PHE:HD1	1:A:492:LEU:H	1.64	0.45
1:B:608:VAL:HA	1:B:692:ILE:CD1	2.47	0.45
1:B:1093:GLY:H	1:B:1107:ARG:NH1	2.15	0.45
1:C:565:PHE:HA	1:C:576:VAL:HA	1.99	0.45
1:C:611:LEU:HB2	1:C:650:LEU:HD13	1.98	0.45
1:C:639:GLY:HA2	1:C:642:VAL:HG23	1.99	0.45
3:E:51:GLY:HA2	3:E:91:TYR:CE2	2.52	0.45
1:A:669:GLY:O	1:A:697:MET:HG2	2.16	0.45
1:B:97:LYS:HG2	1:B:98:SER:H	1.82	0.45
1:B:302:THR:HG21	1:B:314:GLN:O	2.17	0.45
1:C:566:GLY:HA3	1:C:575:ALA:HB3	1.99	0.45
1:C:716:THR:HA	1:C:1110:TYR:HB3	1.99	0.45
1:C:497:PHE:CE2	1:C:507:PRO:HB3	2.52	0.44
2:D:37:VAL:HG21	2:D:134:LEU:CD1	2.47	0.44
1:A:66:HIS:HD2	1:A:263:ALA:O	2.00	0.44
1:A:826:VAL:HB	1:A:1057:PRO:HG2	1.98	0.44
1:B:620:VAL:HG11	1:B:651:ILE:HD12	1.99	0.44
1:A:457:ARG:HD3	1:A:459:SER:O	2.18	0.44
1:A:589:PRO:HB3	1:B:855:PHE:HB3	1.99	0.44
1:A:854:LYS:C	1:A:856:ASN:N	2.74	0.44
1:A:984:LEU:HA	1:C:383:SER:HB3	2.00	0.44
1:B:295:PRO:HG3	1:B:633:TRP:CZ3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:647:ALA:CA	1:C:862:PRO:HG3	2.48	0.44
1:C:152:TRP:CD1	1:C:180:GLU:HB3	2.52	0.44
1:C:171:VAL:HG21	4:b:2:NAG:H81	2.00	0.44
1:C:314:GLN:NE2	1:C:613:GLN:HE22	2.16	0.44
1:C:322:PRO:HG2	1:C:540:ASN:HD21	1.83	0.44
1:A:121:ASN:HA	1:A:126:VAL:HA	1.98	0.44
1:A:615:VAL:H	1:A:648:GLY:HA2	1.82	0.44
1:B:303:LEU:HD11	1:B:313:TYR:CE1	2.52	0.44
1:C:20:THR:HG21	1:C:257:GLY:H	1.82	0.44
1:C:655:HIS:HA	1:C:694:ALA:O	2.18	0.44
1:B:900:MET:HE3	1:B:900:MET:HB2	1.77	0.44
1:B:996:LEU:O	1:B:1000:ARG:HG3	2.17	0.44
2:D:129:GLY:C	2:D:131:GLY:N	2.76	0.44
1:B:448:ASN:ND2	1:B:450:ASN:OD1	2.51	0.44
1:B:571:ASP:HB2	1:C:967:SER:HB3	2.00	0.44
1:B:699:LEU:CD1	1:C:872:GLN:HB2	2.48	0.44
1:C:985:ASP:O	1:C:986:LYS:C	2.58	0.44
2:D:21:VAL:HA	2:D:44:SER:O	2.17	0.44
5:V:1:NAG:H61	5:V:2:NAG:C7	2.47	0.44
1:B:97:LYS:HG2	1:B:98:SER:N	2.32	0.44
1:C:611:LEU:HA	1:C:650:LEU:HA	2.00	0.44
1:C:612:TYR:N	1:C:649:CYS:O	2.24	0.44
1:C:646:ARG:HB2	1:C:668:ALA:CB	2.48	0.44
1:C:852:ALA:O	1:C:855:PHE:HB2	2.17	0.44
1:B:915:VAL:CG2	1:B:1106:GLN:HE22	2.30	0.44
1:C:569:ILE:C	1:C:571:ASP:H	2.25	0.44
1:C:620:VAL:HG21	1:C:651:ILE:HD11	1.99	0.44
1:C:1103:PHE:HZ	6:h:1:NAG:H62	1.83	0.44
1:A:565:PHE:CE1	1:B:42:VAL:HG22	2.52	0.44
1:A:601:GLY:O	1:A:602:THR:C	2.60	0.44
1:A:854:LYS:C	1:A:856:ASN:H	2.26	0.44
1:A:962:LEU:HD11	1:A:1004:LEU:HD23	1.99	0.44
1:A:1101:HIS:CG	6:M:1:NAG:H5	2.52	0.44
1:C:1036:GLN:HG3	1:C:1049:LEU:O	2.18	0.44
4:F:1:NAG:H62	4:F:2:NAG:C7	2.48	0.44
1:A:1030:SER:HB3	1:C:1041:ASP:HB2	1.99	0.43
3:E:81:ARG:NH2	3:E:102:ASP:OD2	2.46	0.43
1:A:186:PHE:O	1:A:211:ASN:HA	2.18	0.43
1:A:299:THR:HG21	1:A:597:VAL:HB	2.00	0.43
1:A:600:PRO:HB2	1:A:604:THR:OG1	2.17	0.43
1:B:1054:GLN:N	1:B:1061:VAL:O	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:599:THR:C	1:C:601:GLY:N	2.77	0.43
2:D:125:VAL:HG21	3:E:109:LEU:CD2	2.49	0.43
1:A:403:ARG:CG	1:A:495:TYR:CE1	2.95	0.43
1:A:1102:TRP:CZ2	1:A:1133:VAL:HG21	2.52	0.43
1:B:133:PHE:CD2	1:B:163:ALA:HB2	2.54	0.43
1:B:395:VAL:HG22	1:B:515:PHE:HB3	2.00	0.43
1:B:759:PHE:HD1	1:B:759:PHE:HA	1.70	0.43
1:C:32:PHE:HB3	1:C:59:PHE:CE2	2.54	0.43
1:C:339:GLY:HA2	4:j:1:NAG:H82	1.98	0.43
6:h:1:NAG:H61	6:h:2:NAG:H82	2.00	0.43
1:B:125:ASN:OD1	4:Q:1:NAG:H5	2.18	0.43
1:C:598:ILE:CD1	1:C:666:ILE:HD11	2.48	0.43
1:C:90:VAL:HG22	1:C:269:TYR:CE2	2.54	0.43
1:C:583:GLU:O	1:C:585:LEU:HG	2.19	0.43
1:B:1107:ARG:HB3	1:C:904:TYR:HE2	1.83	0.43
2:D:37:VAL:HG11	2:D:134:LEU:HD11	2.00	0.43
1:A:736:VAL:HG13	1:A:858:LEU:HD23	2.01	0.43
2:D:37:VAL:HG22	2:D:38:ARG:H	1.84	0.43
2:D:109:ASP:O	2:D:136:VAL:HG23	2.19	0.43
1:A:1062:PHE:HB2	1:A:1064:HIS:NE2	2.34	0.43
1:B:102:ARG:HB3	1:B:121:ASN:O	2.18	0.43
1:A:609:ALA:HB2	1:A:692:ILE:HG21	2.00	0.43
1:A:979:ASP:OD1	1:C:518:LEU:HD13	2.19	0.43
1:C:196:ASN:HA	1:C:201:PHE:HA	2.01	0.43
1:A:394:ASN:HA	1:A:524:VAL:HG21	2.01	0.42
1:A:819:GLU:HG3	1:A:1054:GLN:OE1	2.19	0.42
1:C:474:GLN:HG2	1:C:480:CYS:HB3	2.01	0.42
1:C:909:ILE:HD13	1:C:1049:LEU:HD21	2.01	0.42
2:D:66:TRP:CZ2	2:D:68:GLY:HA2	2.54	0.42
1:B:124:THR:CB	4:Q:1:NAG:HN2	2.32	0.42
1:B:722:VAL:HA	1:B:1064:HIS:O	2.19	0.42
1:B:1077:THR:HG23	1:C:900:MET:HE3	2.00	0.42
1:A:909:ILE:HD13	1:A:1049:LEU:HD21	2.01	0.42
1:B:567:ARG:CZ	1:C:42:VAL:HG21	2.48	0.42
1:C:106:PHE:HB3	1:C:235:ILE:HG23	2.01	0.42
1:C:141:LEU:O	1:C:243:ALA:HA	2.19	0.42
1:C:520:ALA:HB1	1:C:521:PRO:HD2	2.01	0.42
1:C:559:PHE:HB2	1:C:584:ILE:HG13	2.00	0.42
1:A:930:ALA:O	1:A:934:ILE:HG12	2.19	0.42
1:B:1054:GLN:HB2	1:B:1061:VAL:HB	2.01	0.42
1:C:110:LEU:HB3	1:C:135:PHE:HD2	1.80	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:643:PHE:CE2	1:C:645:THR:HG22	2.54	0.42
1:A:439:ASN:HB2	1:A:506:GLN:HB3	2.02	0.42
1:B:609:ALA:HB1	1:B:653:ALA:HB2	2.01	0.42
1:B:699:LEU:HD12	1:C:872:GLN:HB3	2.02	0.42
1:C:121:ASN:CG	1:C:126:VAL:HG22	2.44	0.42
1:A:299:THR:CG2	1:A:597:VAL:HB	2.50	0.42
1:C:497:PHE:CD2	1:C:507:PRO:HB3	2.55	0.42
1:C:553:THR:O	1:C:586:ASP:N	2.52	0.42
1:A:715:PRO:HD3	1:B:894:LEU:HD13	2.01	0.42
1:B:383:SER:HB2	1:C:985:ASP:N	2.34	0.42
1:B:984:LEU:HD23	1:B:988:GLU:HB3	2.01	0.42
1:B:1017:GLU:HG2	1:C:1019:ARG:HH22	1.85	0.42
1:C:562:PHE:CE2	1:C:563:GLN:HG3	2.54	0.42
1:C:736:VAL:HG13	1:C:858:LEU:HD23	2.01	0.42
1:A:987:VAL:HG21	1:B:413:GLY:HA3	2.02	0.42
1:A:1148:PHE:HE1	1:C:1152:LEU:HD13	1.84	0.42
1:C:1086:LYS:HA	1:C:1125:ASN:HA	2.02	0.42
1:A:290:ASP:HB3	1:A:293:LEU:HB2	2.02	0.42
1:A:352:ALA:CB	1:A:466:ARG:HH21	2.23	0.42
1:A:702:GLU:HG2	1:B:790:LYS:HD2	2.02	0.42
1:B:97:LYS:C	1:B:99:ASN:H	2.28	0.42
1:C:105:ILE:HG23	1:C:135:PHE:CZ	2.54	0.42
1:A:901:GLN:O	1:A:905:ARG:HG2	2.20	0.41
1:B:390:LEU:CD2	1:C:983:ARG:HG2	2.45	0.41
1:A:852:ALA:HB1	1:C:570:ALA:CB	2.47	0.41
1:A:853:GLN:HG2	1:A:963:VAL:HG21	2.02	0.41
1:B:618:THR:HG21	4:S:1:NAG:H5	2.02	0.41
3:E:36:GLY:H	3:E:98:PRO:HD2	1.85	0.41
1:A:894:LEU:HD13	1:C:715:PRO:HD3	2.02	0.41
1:B:1048:HIS:HA	1:B:1066:THR:HG22	2.00	0.41
1:A:287:ASP:HB3	1:A:306:PHE:HE2	1.85	0.41
1:A:476:GLY:N	1:A:487:ASN:HB3	2.36	0.41
1:A:665:PRO:CB	1:B:864:LEU:HD22	2.50	0.41
1:A:815:ARG:HG2	1:A:871:ALA:CB	2.50	0.41
1:B:1148:PHE:CD2	1:C:1148:PHE:CE1	3.08	0.41
1:C:611:LEU:HD12	1:C:650:LEU:HB2	2.03	0.41
1:C:914:ASN:O	1:C:915:VAL:C	2.62	0.41
1:A:431:GLY:HA2	1:A:515:PHE:HD2	1.85	0.41
1:B:456:PHE:HB3	1:B:473:TYR:CD2	2.55	0.41
1:B:995:ARG:NH1	1:C:994:ASP:OD1	2.52	0.41
1:C:83:VAL:HA	1:C:239:GLN:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:ASP:C	1:C:200:TYR:H	2.29	0.41
2:D:102:LEU:HB3	2:D:105:LEU:CD2	2.29	0.41
2:D:102:LEU:HD13	2:D:134:LEU:CD2	2.50	0.41
1:A:106:PHE:HB2	1:A:117:LEU:HB3	2.02	0.41
1:A:560:LEU:O	1:A:562:PHE:N	2.54	0.41
1:A:563:GLN:NE2	1:B:43:PHE:HD2	2.17	0.41
1:B:193:VAL:HG13	1:B:270:LEU:HD11	2.03	0.41
1:B:214:ARG:O	1:B:214:ARG:HG3	2.21	0.41
1:B:665:PRO:HB2	1:C:864:LEU:HD22	2.03	0.41
1:B:826:VAL:HB	1:B:1057:PRO:HG2	2.02	0.41
1:B:1009:THR:HA	1:B:1012:LEU:HD12	2.02	0.41
1:B:1017:GLU:HG2	1:C:1019:ARG:NH2	2.35	0.41
1:B:1070:ALA:O	1:B:1071:GLN:C	2.62	0.41
1:C:581:THR:O	1:C:583:GLU:HG2	2.21	0.41
1:B:339:GLY:HA2	4:Y:1:NAG:C8	2.48	0.41
2:D:23:LEU:HD22	2:D:129:GLY:N	2.35	0.41
1:A:83:VAL:HG11	1:A:237:ARG:NH2	2.34	0.41
1:A:206:LYS:HB3	1:A:223:LEU:CD2	2.51	0.41
1:B:571:ASP:HB2	1:C:967:SER:CB	2.51	0.41
1:C:121:ASN:C	1:C:121:ASN:HD22	2.28	0.41
1:C:319:ARG:CZ	1:C:319:ARG:CB	2.99	0.41
1:C:636:TYR:HD2	1:C:651:ILE:HG22	1.86	0.41
2:D:39:ILE:CD1	2:D:134:LEU:HD22	2.51	0.41
1:A:42:VAL:HA	1:C:565:PHE:O	2.21	0.41
1:A:200:TYR:OH	1:C:394:ASN:HB2	2.16	0.41
1:B:392:PHE:CD2	1:B:517:LEU:HD21	2.56	0.41
1:B:779:GLN:O	1:B:780:GLU:C	2.64	0.41
1:C:49:HIS:HB3	1:C:279:TYR:CE1	2.53	0.41
1:C:66:HIS:N	1:C:80:ASP:OD2	2.54	0.41
1:C:805:ILE:HB	1:C:1054:GLN:NE2	2.36	0.41
1:B:105:ILE:CD1	1:B:118:LEU:HD12	2.51	0.40
1:B:396:TYR:OH	1:C:230:PRO:HB3	2.21	0.40
1:B:1126:CYS:O	1:B:1127:ASP:C	2.64	0.40
1:A:188:ASN:HB2	1:A:207:HIS:NE2	2.35	0.40
1:A:369:TYR:OH	1:C:475:ALA:HB1	2.22	0.40
1:A:1142:GLN:O	1:A:1146:ASP:N	2.49	0.40
1:C:196:ASN:HB2	1:C:201:PHE:CD1	2.56	0.40
1:C:322:PRO:HG3	1:C:549:THR:HG23	2.04	0.40
1:C:328:ARG:HD3	1:C:533:LEU:HB2	2.03	0.40
1:C:335:LEU:C	1:C:361:CYS:HB2	2.46	0.40
1:A:470:THR:HB	1:A:490:PHE:CZ	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:609:ALA:HB3	1:A:692:ILE:HG21	2.02	0.40
1:A:864:LEU:HA	1:C:667:GLY:HA2	2.04	0.40
1:A:982:SER:HB2	1:C:390:LEU:HD11	2.04	0.40
1:C:32:PHE:O	1:C:33:THR:C	2.64	0.40
1:A:200:TYR:HH	1:C:394:ASN:CB	2.34	0.40
1:B:1123:SER:OG	1:C:918:GLU:OE2	2.39	0.40
1:C:653:ALA:HA	1:C:692:ILE:O	2.20	0.40
1:A:527:PRO:O	1:A:527:PRO:CG	2.70	0.40
1:B:189:LEU:HD22	1:B:217:PRO:CG	2.50	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	952/1135 (84%)	944 (99%)	8 (1%)	73	90
1	B	952/1135 (84%)	947 (100%)	5 (0%)	81	93
1	C	961/1135 (85%)	949 (99%)	12 (1%)	63	87
2	D	105/120 (88%)	101 (96%)	4 (4%)	29	64
3	E	91/109 (84%)	90 (99%)	1 (1%)	65	87
All	All	3061/3634 (84%)	3031 (99%)	30 (1%)	65	88

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	LEU
1	A	125	ASN
1	A	177	MET
1	A	179	LEU
1	A	492	LEU
1	A	493	GLN
1	A	565	PHE
1	A	878	LEU
1	B	66	HIS
1	B	118	LEU
1	B	568	ASP
1	B	759	PHE
1	B	916	LEU
1	C	18	LEU
1	C	20	THR
1	C	32	PHE
1	C	80	ASP
1	C	108	THR
1	C	212	LEU
1	C	226	LEU
1	C	625	HIS
1	C	825	LYS
1	C	1038	LYS
1	C	1050	MET
1	C	1086	LYS
2	D	20	GLU
2	D	24	GLN
2	D	116	THR
2	D	125	VAL
3	E	74	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (74) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	99	ASN
1	A	115	GLN
1	A	125	ASN
1	A	149	ASN
1	A	271	GLN
1	A	314	GLN
1	A	394	ASN
1	A	460	ASN

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Mol	Chain	Res	Type
1	A	493	GLN
1	A	501	ASN
1	A	540	ASN
1	A	542	ASN
1	A	563	GLN
1	A	580	GLN
1	A	606	ASN
1	A	644	GLN
1	A	675	GLN
1	A	690	GLN
1	A	703	ASN
1	A	764	ASN
1	A	804	GLN
1	A	913	GLN
1	A	935	GLN
1	A	965	GLN
1	A	969	ASN
1	A	1010	GLN
1	A	1083	HIS
1	A	1108	ASN
1	A	1159	HIS
1	B	87	ASN
1	B	149	ASN
1	B	314	GLN
1	B	360	ASN
1	B	448	ASN
1	B	658	ASN
1	B	779	GLN
1	B	787	GLN
1	B	804	GLN
1	B	913	GLN
1	B	960	ASN
1	B	978	ASN
1	B	1002	GLN
1	B	1106	GLN
1	B	1108	ASN
1	B	1142	GLN
1	C	52	GLN
1	C	115	GLN
1	C	134	GLN
1	C	149	ASN
1	C	207	HIS

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Mol	Chain	Res	Type
1	C	211	ASN
1	C	314	GLN
1	C	394	ASN
1	C	474	GLN
1	C	498	GLN
1	C	540	ASN
1	C	544	ASN
1	C	580	GLN
1	C	607	GLN
1	C	644	GLN
1	C	690	GLN
1	C	755	GLN
1	C	779	GLN
1	C	804	GLN
1	C	907	ASN
1	C	913	GLN
1	C	935	GLN
1	C	949	GLN
1	C	953	ASN
1	C	969	ASN
1	C	1002	GLN
1	C	1005	GLN
1	C	1135	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

72 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	$\# Z > 2$
4	NAG	F	1	4,1	14,14,15	0.21	0	17,19,21	0.47	0
4	NAG	F	2	4	14,14,15	0.27	0	17,19,21	0.55	0
4	NAG	G	1	4,1	14,14,15	0.40	0	17,19,21	0.45	0
4	NAG	G	2	4	14,14,15	0.39	0	17,19,21	0.38	0
4	NAG	H	1	4,1	14,14,15	0.24	0	17,19,21	0.50	0
4	NAG	H	2	4	14,14,15	0.29	0	17,19,21	0.50	0
4	NAG	I	1	4,1	14,14,15	0.40	0	17,19,21	0.35	0
4	NAG	I	2	4	14,14,15	0.41	0	17,19,21	0.35	0
4	NAG	J	1	4,1	14,14,15	0.39	0	17,19,21	0.62	1 (5%)
4	NAG	J	2	4	14,14,15	0.40	0	17,19,21	0.41	0
4	NAG	K	1	4,1	14,14,15	0.38	0	17,19,21	0.44	0
4	NAG	K	2	4	14,14,15	0.40	0	17,19,21	0.40	0
5	NAG	L	1	5,1	14,14,15	0.41	0	17,19,21	0.68	0
5	NAG	L	2	5	14,14,15	0.40	0	17,19,21	0.44	0
5	FUC	L	3	5	10,10,11	1.65	2 (20%)	14,14,16	0.90	0
6	NAG	M	1	6,1	14,14,15	0.40	0	17,19,21	0.48	0
6	NAG	M	2	6	14,14,15	0.39	0	17,19,21	0.38	0
6	MAN	M	3	6	11,11,12	0.22	0	15,15,17	0.44	0
6	NAG	N	1	6,1	14,14,15	0.39	0	17,19,21	0.40	0
6	NAG	N	2	6	14,14,15	0.40	0	17,19,21	0.40	0
6	MAN	N	3	6	11,11,12	0.20	0	15,15,17	0.62	0
4	NAG	O	1	4,1	14,14,15	0.39	0	17,19,21	0.40	0
4	NAG	O	2	4	14,14,15	0.40	0	17,19,21	0.35	0
4	NAG	P	1	4,1	14,14,15	0.40	0	17,19,21	0.63	0
4	NAG	P	2	4	14,14,15	0.30	0	17,19,21	0.53	0
4	NAG	Q	1	4,1	14,14,15	0.39	0	17,19,21	0.67	1 (5%)
4	NAG	Q	2	4	14,14,15	0.39	0	17,19,21	0.37	0
4	NAG	R	1	4,1	14,14,15	0.39	0	17,19,21	0.45	0
4	NAG	R	2	4	14,14,15	0.40	0	17,19,21	0.37	0
4	NAG	S	1	4,1	14,14,15	0.33	0	17,19,21	0.43	0
4	NAG	S	2	4	14,14,15	0.29	0	17,19,21	0.54	0
4	NAG	T	1	4,1	14,14,15	0.38	0	17,19,21	0.67	1 (5%)
4	NAG	T	2	4	14,14,15	0.39	0	17,19,21	0.41	0
4	NAG	U	1	4,1	14,14,15	0.40	0	17,19,21	0.43	0
4	NAG	U	2	4	14,14,15	0.39	0	17,19,21	0.38	0
5	NAG	V	1	5,1	14,14,15	0.44	0	17,19,21	0.58	0
5	NAG	V	2	5	14,14,15	0.39	0	17,19,21	0.48	0
5	FUC	V	3	5	10,10,11	1.58	2 (20%)	14,14,16	1.07	1 (7%)
6	NAG	W	1	6,1	14,14,15	0.39	0	17,19,21	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	W	2	6	14,14,15	0.40	0	17,19,21	0.35	0
6	MAN	W	3	6	11,11,12	0.18	0	15,15,17	0.47	0
6	NAG	X	1	6,1	14,14,15	0.39	0	17,19,21	0.50	0
6	NAG	X	2	6	14,14,15	0.39	0	17,19,21	0.37	0
6	MAN	X	3	6	11,11,12	0.20	0	15,15,17	0.60	0
4	NAG	Y	1	4,1	14,14,15	0.73	1 (7%)	17,19,21	0.50	0
4	NAG	Y	2	4	14,14,15	0.27	0	17,19,21	0.57	0
4	NAG	Z	1	4,1	14,14,15	0.39	0	17,19,21	0.39	0
4	NAG	Z	2	4	14,14,15	0.40	0	17,19,21	0.37	0
4	NAG	a	1	4,1	14,14,15	0.24	0	17,19,21	0.50	0
4	NAG	a	2	4	14,14,15	0.30	0	17,19,21	0.54	0
4	NAG	b	1	4,1	14,14,15	0.40	0	17,19,21	0.29	0
4	NAG	b	2	4	14,14,15	0.39	0	17,19,21	0.43	0
4	NAG	c	1	4,1	14,14,15	0.39	0	17,19,21	0.57	0
4	NAG	c	2	4	14,14,15	0.39	0	17,19,21	0.44	0
4	NAG	d	1	4,1	14,14,15	0.40	0	17,19,21	0.41	0
4	NAG	d	2	4	14,14,15	0.39	0	17,19,21	0.37	0
4	NAG	e	1	4,1	14,14,15	0.39	0	17,19,21	0.60	1 (5%)
4	NAG	e	2	4	14,14,15	0.39	0	17,19,21	0.41	0
4	NAG	f	1	4,1	14,14,15	0.38	0	17,19,21	0.38	0
4	NAG	f	2	4	14,14,15	0.40	0	17,19,21	0.37	0
5	NAG	g	1	5,1	14,14,15	0.41	0	17,19,21	0.87	2 (11%)
5	NAG	g	2	5	14,14,15	0.39	0	17,19,21	0.40	0
5	FUC	g	3	5	10,10,11	1.68	2 (20%)	14,14,16	1.13	1 (7%)
6	NAG	h	1	6,1	14,14,15	0.40	0	17,19,21	0.38	0
6	NAG	h	2	6	14,14,15	0.40	0	17,19,21	0.38	0
6	MAN	h	3	6	11,11,12	0.23	0	15,15,17	0.49	0
4	NAG	i	1	4,1	14,14,15	0.40	0	17,19,21	0.35	0
4	NAG	i	2	4	14,14,15	0.39	0	17,19,21	0.42	0
4	NAG	j	1	4,1	14,14,15	0.55	0	17,19,21	0.50	0
4	NAG	j	2	4	14,14,15	0.28	0	17,19,21	0.55	0
4	NAG	k	1	4,1	14,14,15	0.39	0	17,19,21	0.54	0
4	NAG	k	2	4	14,14,15	0.39	0	17,19,21	0.47	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
4	NAG	F	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1

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

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

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	L	3	FUC	O5-C5	3.10	1.49	1.43
5	g	3	FUC	O5-C5	2.93	1.49	1.43
5	V	3	FUC	O5-C5	2.91	1.49	1.43
5	g	3	FUC	O5-C1	2.64	1.48	1.43
4	Y	1	NAG	O5-C1	-2.63	1.39	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	V	3	FUC	O5-C1	2.26	1.47	1.43
5	L	3	FUC	O5-C1	2.20	1.47	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	g	3	FUC	C1-C2-C3	2.51	113.30	109.64
4	T	1	NAG	O5-C1-C2	-2.36	107.64	111.29
4	Q	1	NAG	O5-C1-C2	-2.23	107.84	111.29
5	g	1	NAG	C2-N2-C7	2.22	125.88	122.90
5	g	1	NAG	C1-C2-N2	2.13	113.79	110.43
4	J	1	NAG	O5-C1-C2	-2.10	108.04	111.29
5	V	3	FUC	C1-C2-C3	2.10	112.70	109.64
4	e	1	NAG	O5-C1-C2	-2.00	108.19	111.29

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	I	1	NAG	C8-C7-N2-C2
4	I	1	NAG	O7-C7-N2-C2
4	K	1	NAG	O7-C7-N2-C2
4	U	1	NAG	O7-C7-N2-C2
4	d	1	NAG	C8-C7-N2-C2
4	d	1	NAG	O7-C7-N2-C2
4	d	2	NAG	C8-C7-N2-C2
4	d	2	NAG	O7-C7-N2-C2
4	f	1	NAG	O7-C7-N2-C2
5	g	1	NAG	C1-C2-N2-C7
6	M	1	NAG	C8-C7-N2-C2
6	M	1	NAG	O7-C7-N2-C2
6	N	1	NAG	C8-C7-N2-C2
6	N	1	NAG	O7-C7-N2-C2
4	K	1	NAG	C8-C7-N2-C2
4	U	1	NAG	C8-C7-N2-C2
4	f	1	NAG	C8-C7-N2-C2
4	Y	1	NAG	O5-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6
5	L	1	NAG	O5-C5-C6-O6
4	I	2	NAG	C8-C7-N2-C2
4	I	2	NAG	O7-C7-N2-C2
6	N	2	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
4	a	1	NAG	O5-C5-C6-O6
4	j	1	NAG	O5-C5-C6-O6
4	Y	1	NAG	C4-C5-C6-O6
5	g	1	NAG	O5-C5-C6-O6
5	V	1	NAG	O5-C5-C6-O6
4	I	1	NAG	C4-C5-C6-O6
4	j	1	NAG	C4-C5-C6-O6
4	k	2	NAG	C8-C7-N2-C2
6	N	2	NAG	C8-C7-N2-C2
5	V	1	NAG	C4-C5-C6-O6
5	g	1	NAG	C4-C5-C6-O6
4	k	2	NAG	O7-C7-N2-C2
4	a	1	NAG	C4-C5-C6-O6
5	L	1	NAG	C4-C5-C6-O6
4	R	2	NAG	C8-C7-N2-C2
4	R	2	NAG	O7-C7-N2-C2
4	d	1	NAG	O5-C5-C6-O6
6	M	2	NAG	C8-C7-N2-C2
4	f	2	NAG	C8-C7-N2-C2
6	M	2	NAG	O7-C7-N2-C2
5	g	1	NAG	C8-C7-N2-C2
4	R	1	NAG	C8-C7-N2-C2
4	Q	2	NAG	C8-C7-N2-C2
4	f	2	NAG	O7-C7-N2-C2
4	c	2	NAG	O5-C5-C6-O6
5	g	1	NAG	O7-C7-N2-C2
6	W	3	MAN	O5-C5-C6-O6
5	g	1	NAG	C3-C2-N2-C7
4	d	1	NAG	C4-C5-C6-O6
4	R	1	NAG	O7-C7-N2-C2
4	Q	2	NAG	O7-C7-N2-C2
4	k	1	NAG	C8-C7-N2-C2
4	k	1	NAG	O7-C7-N2-C2
4	R	1	NAG	C4-C5-C6-O6
4	S	1	NAG	C1-C2-N2-C7
4	J	2	NAG	C3-C2-N2-C7
4	T	2	NAG	C3-C2-N2-C7
4	e	2	NAG	C3-C2-N2-C7
6	M	2	NAG	O5-C5-C6-O6
4	K	1	NAG	C4-C5-C6-O6
6	h	2	NAG	O5-C5-C6-O6

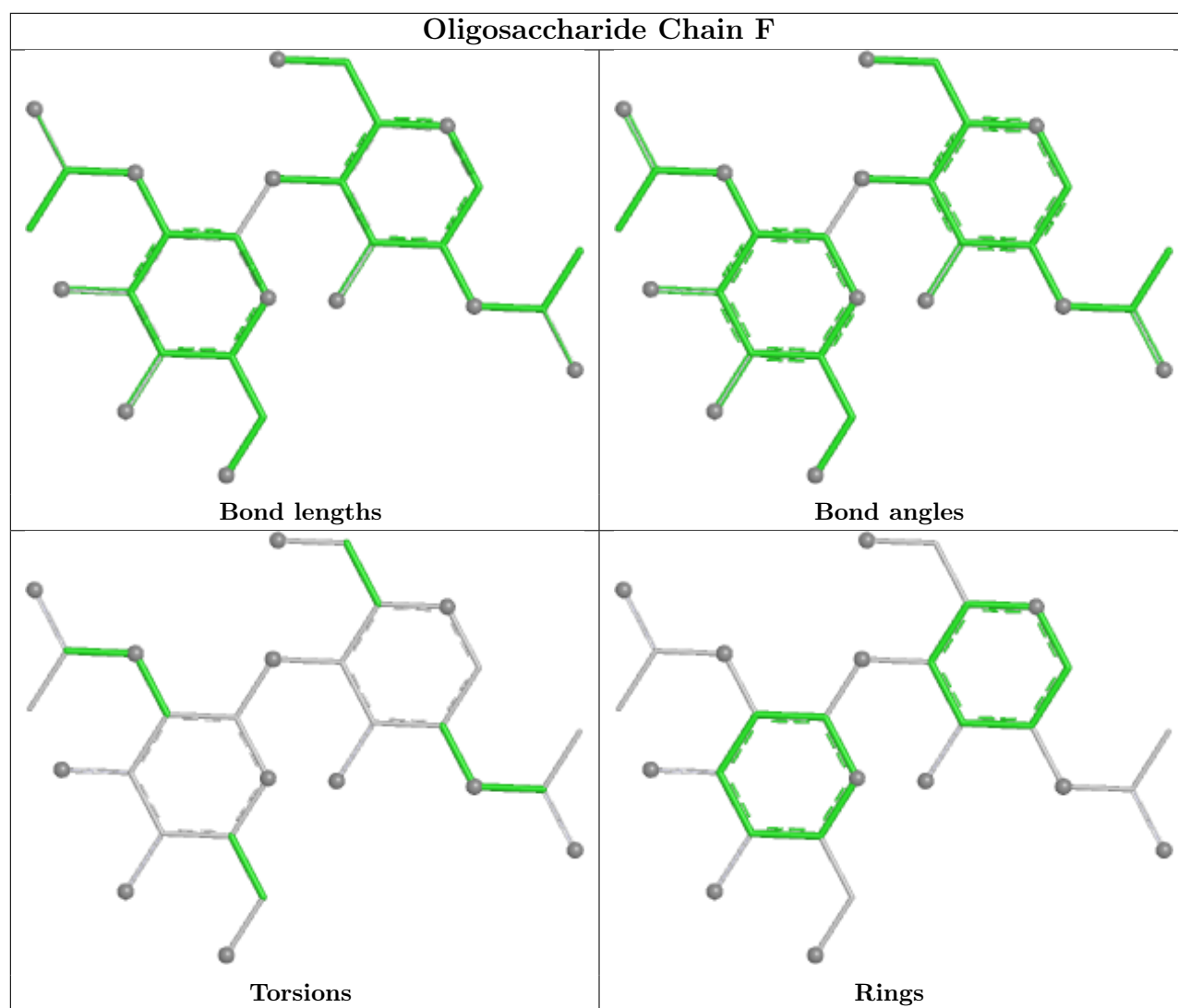
All (3) ring outliers are listed below:

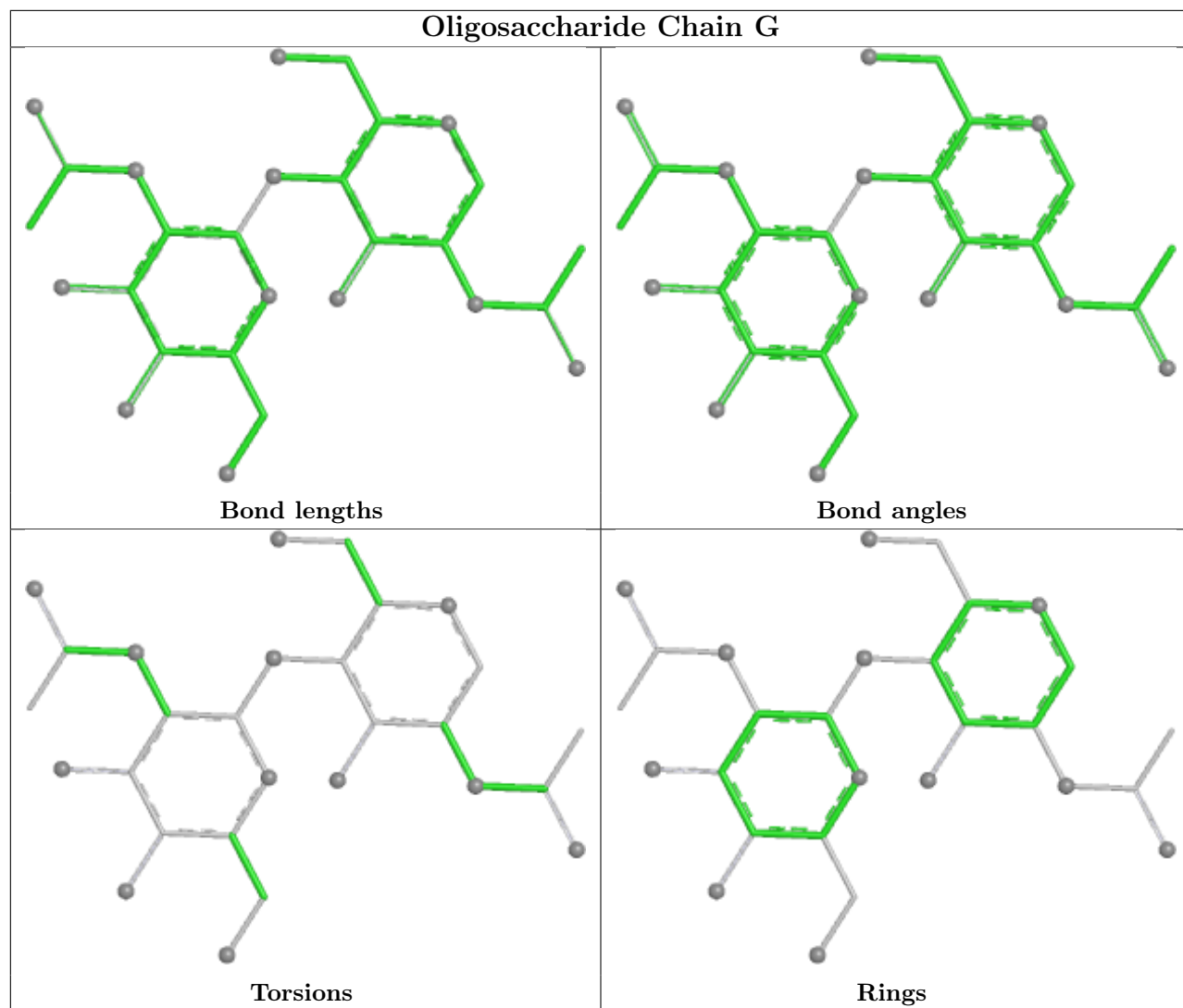
Mol	Chain	Res	Type	Atoms
6	M	3	MAN	C1-C2-C3-C4-C5-O5
6	X	3	MAN	C1-C2-C3-C4-C5-O5
6	N	3	MAN	C1-C2-C3-C4-C5-O5

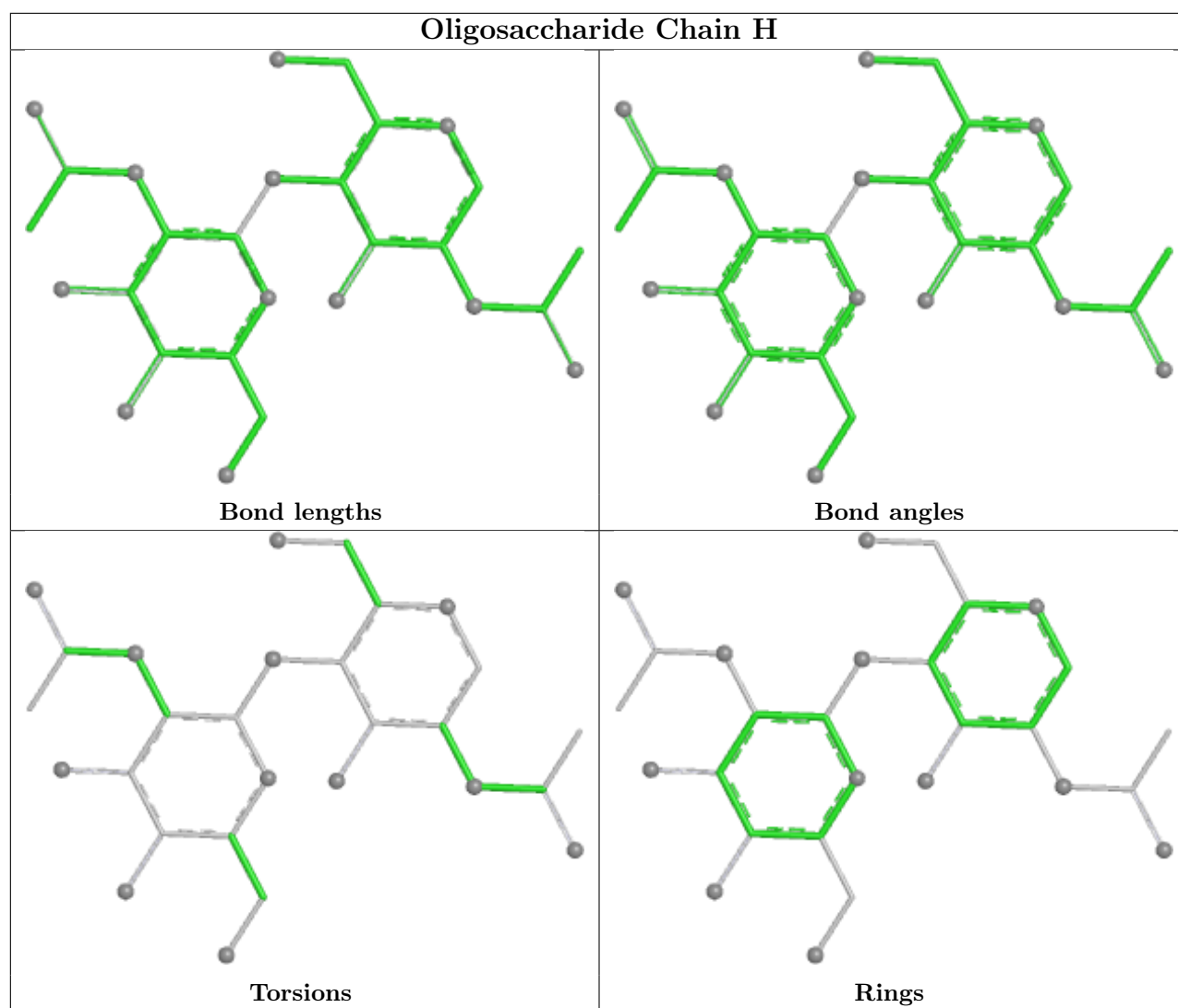
24 monomers are involved in 31 short contacts:

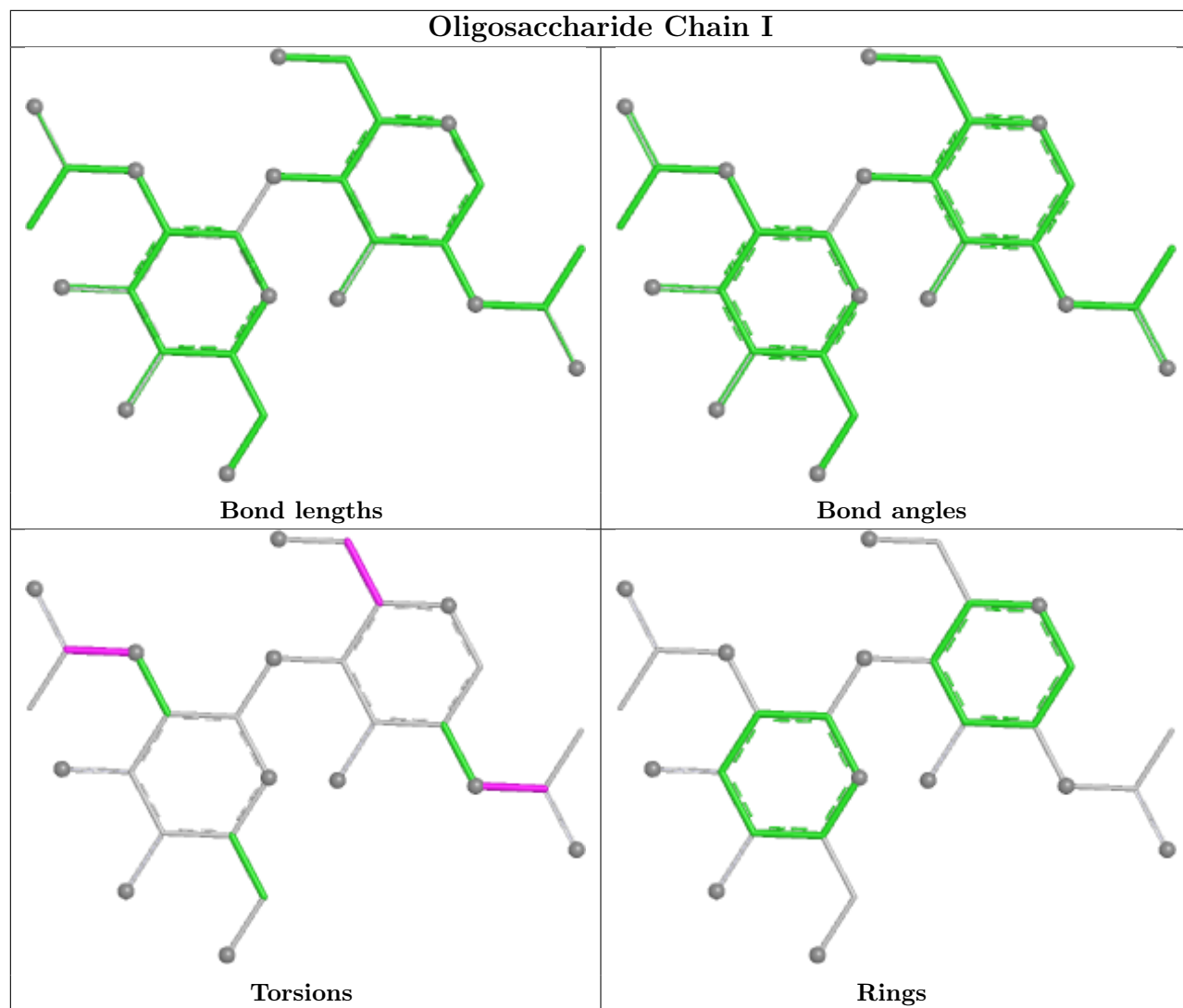
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	h	1	NAG	3	0
6	M	1	NAG	2	0
6	h	2	NAG	1	0
4	e	2	NAG	1	0
4	P	1	NAG	1	0
4	S	1	NAG	1	0
4	G	1	NAG	2	0
4	j	1	NAG	1	0
6	W	1	NAG	1	0
5	g	3	FUC	1	0
6	M	2	NAG	1	0
4	J	2	NAG	1	0
4	Y	1	NAG	3	0
4	F	1	NAG	1	0
5	g	2	NAG	1	0
4	b	2	NAG	1	0
4	k	1	NAG	1	0
4	Q	1	NAG	4	0
4	a	1	NAG	4	0
4	O	1	NAG	1	0
4	F	2	NAG	1	0
5	V	1	NAG	1	0
4	b	1	NAG	1	0
5	V	2	NAG	1	0

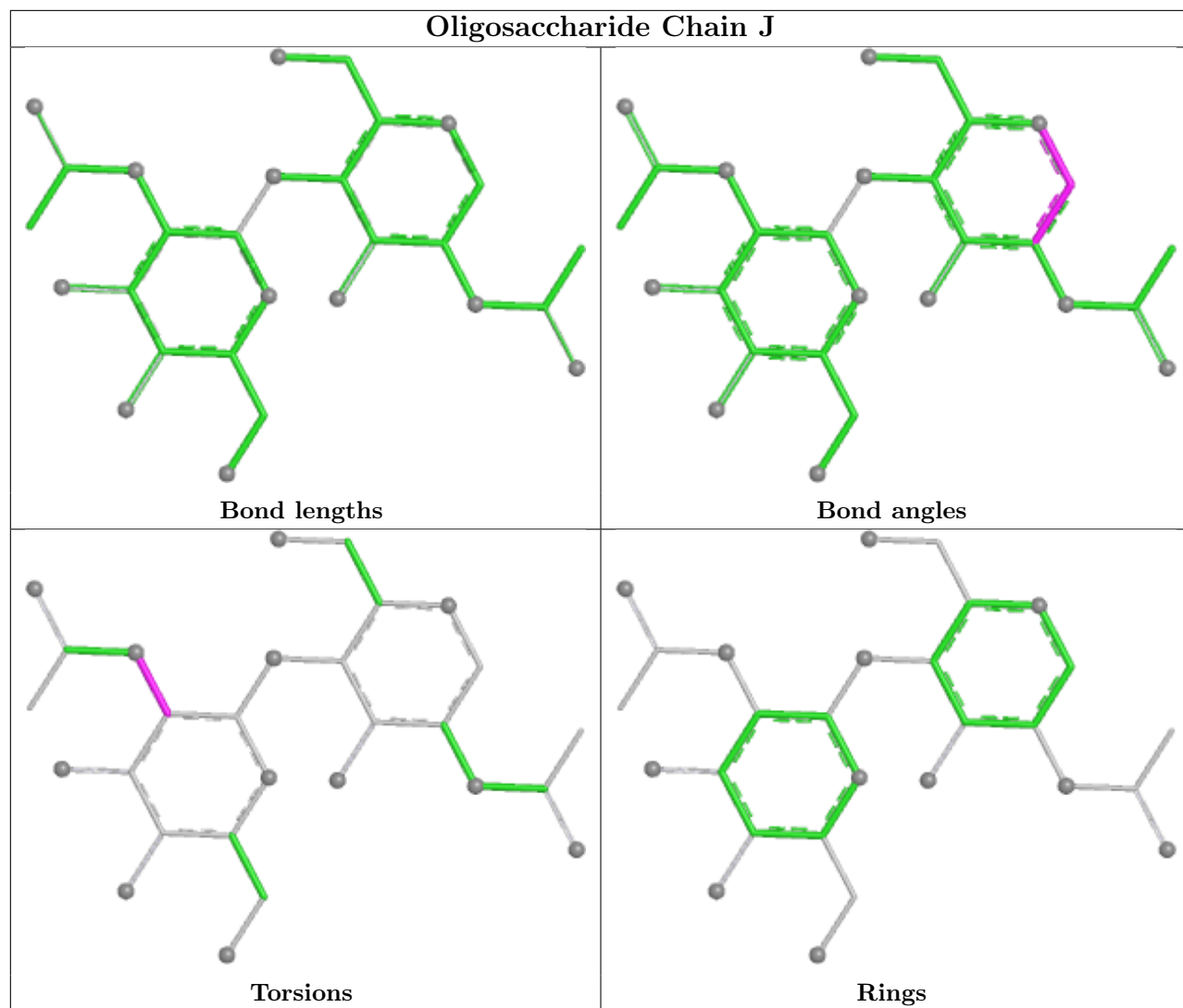
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

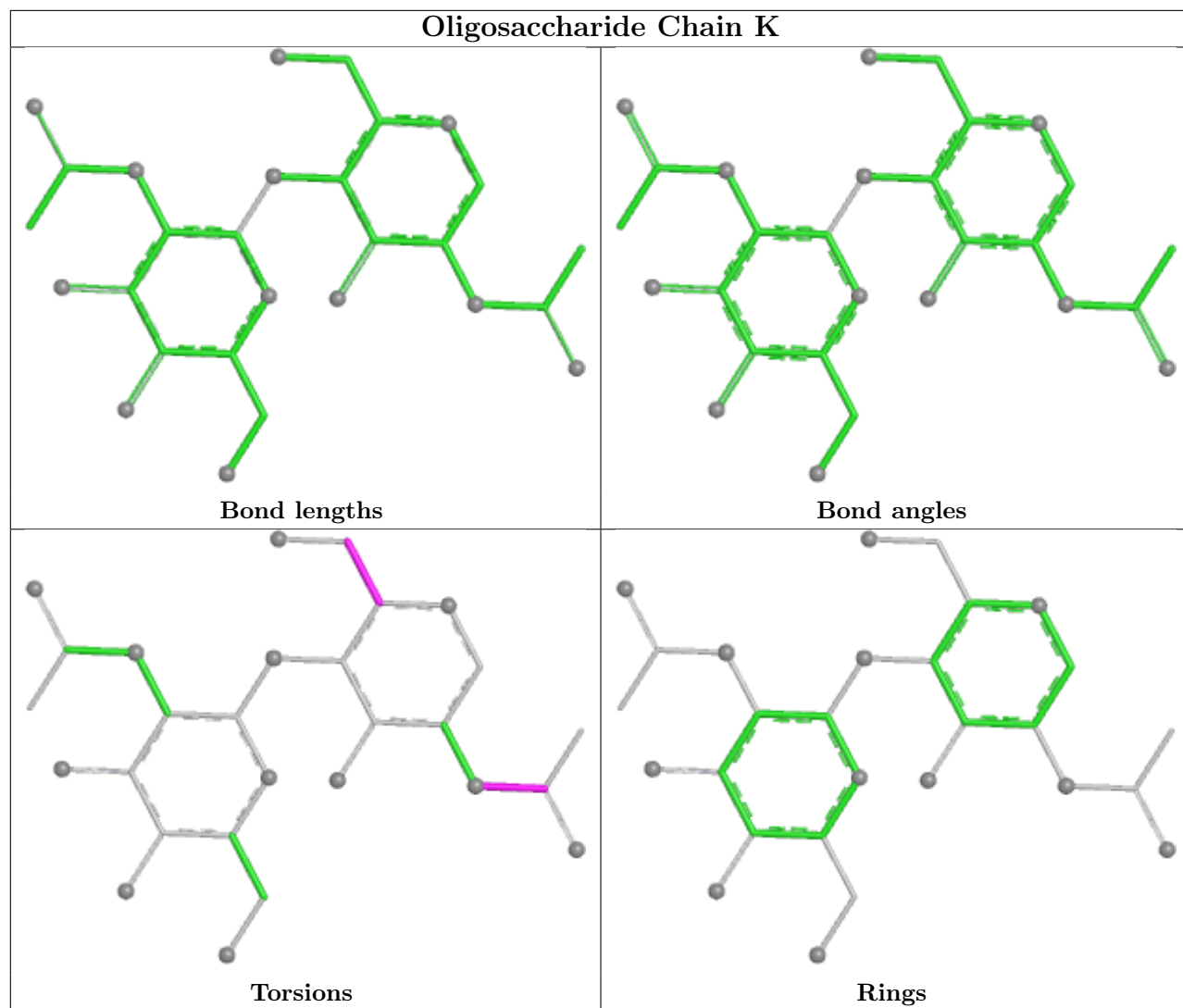


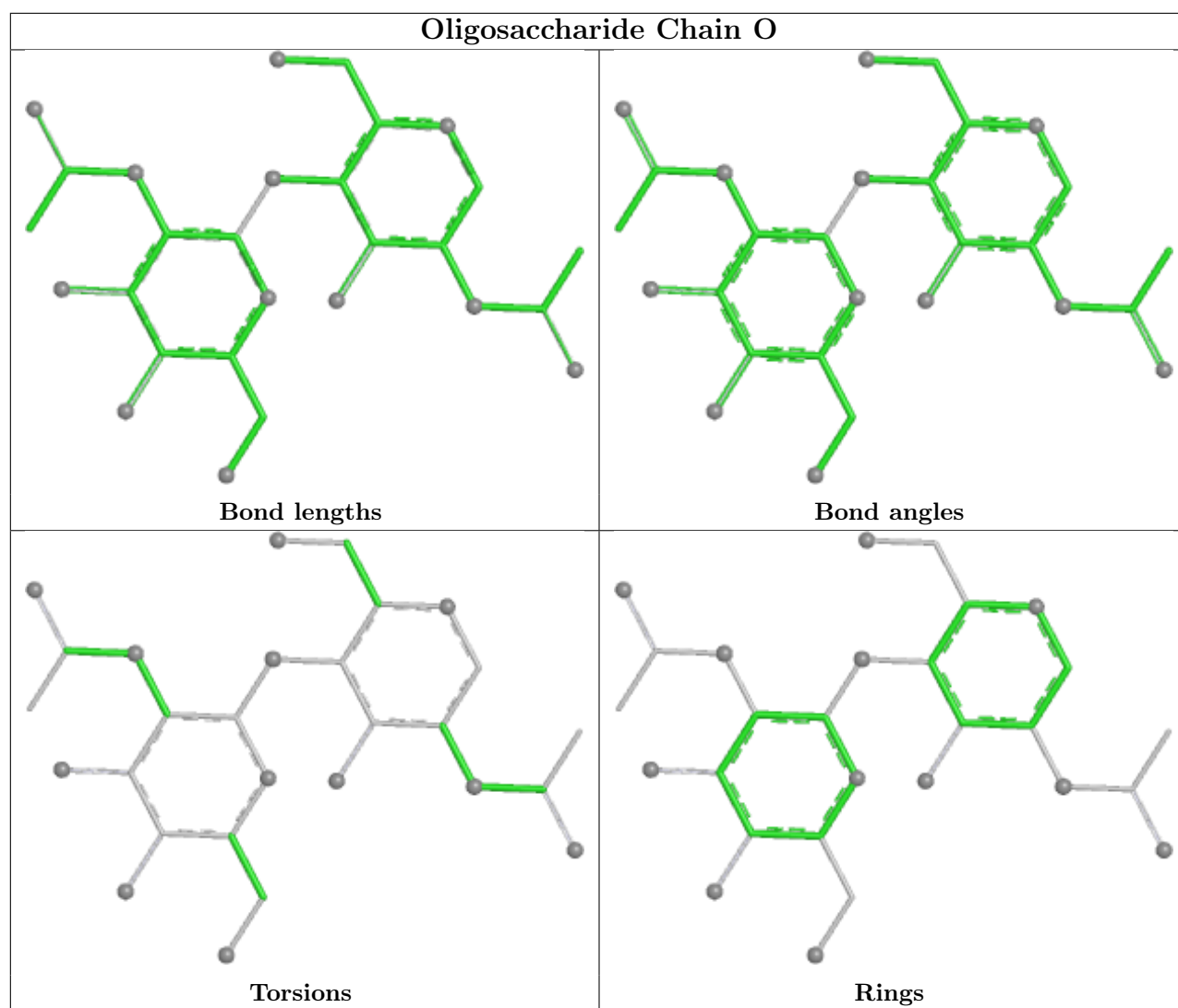


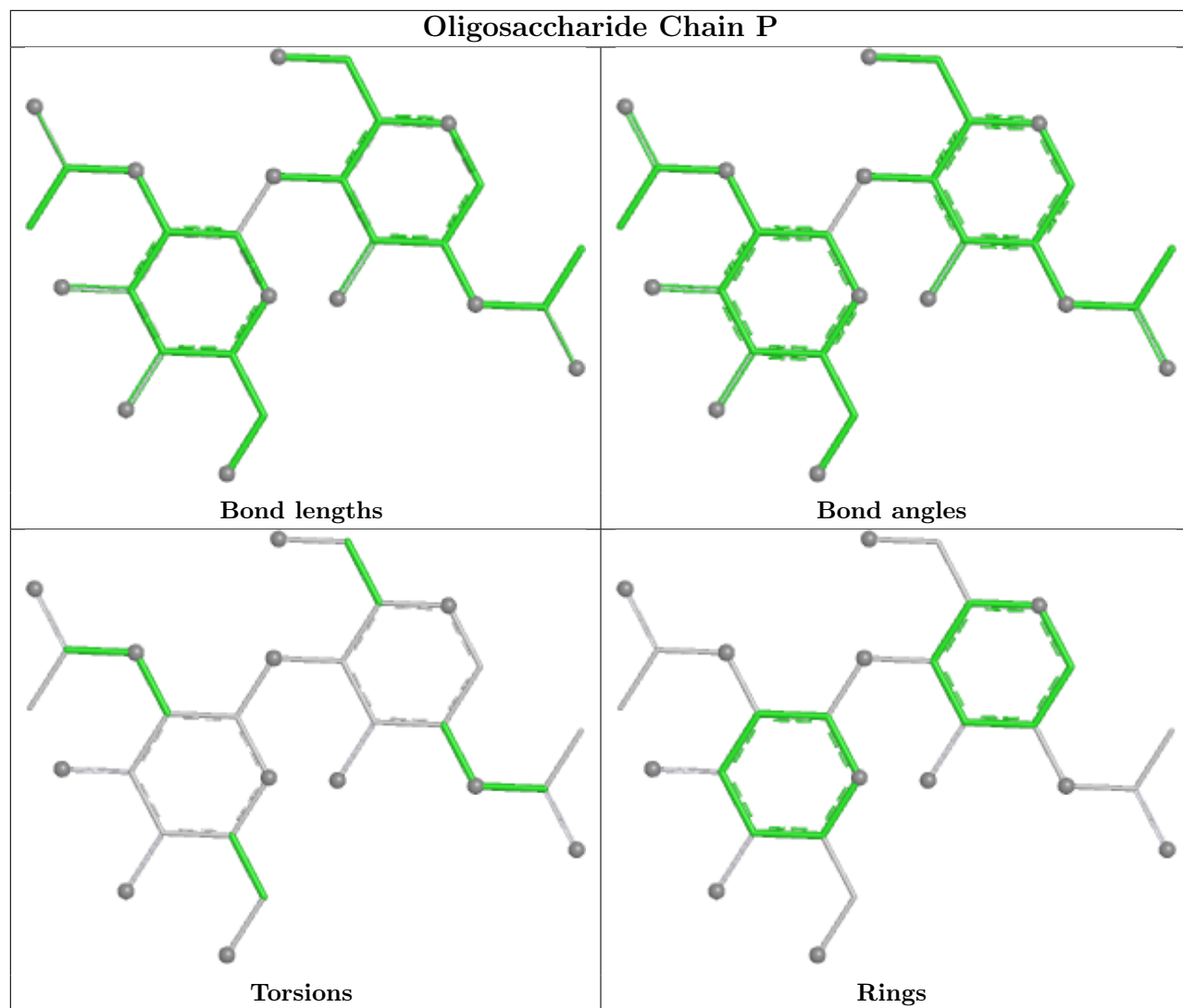


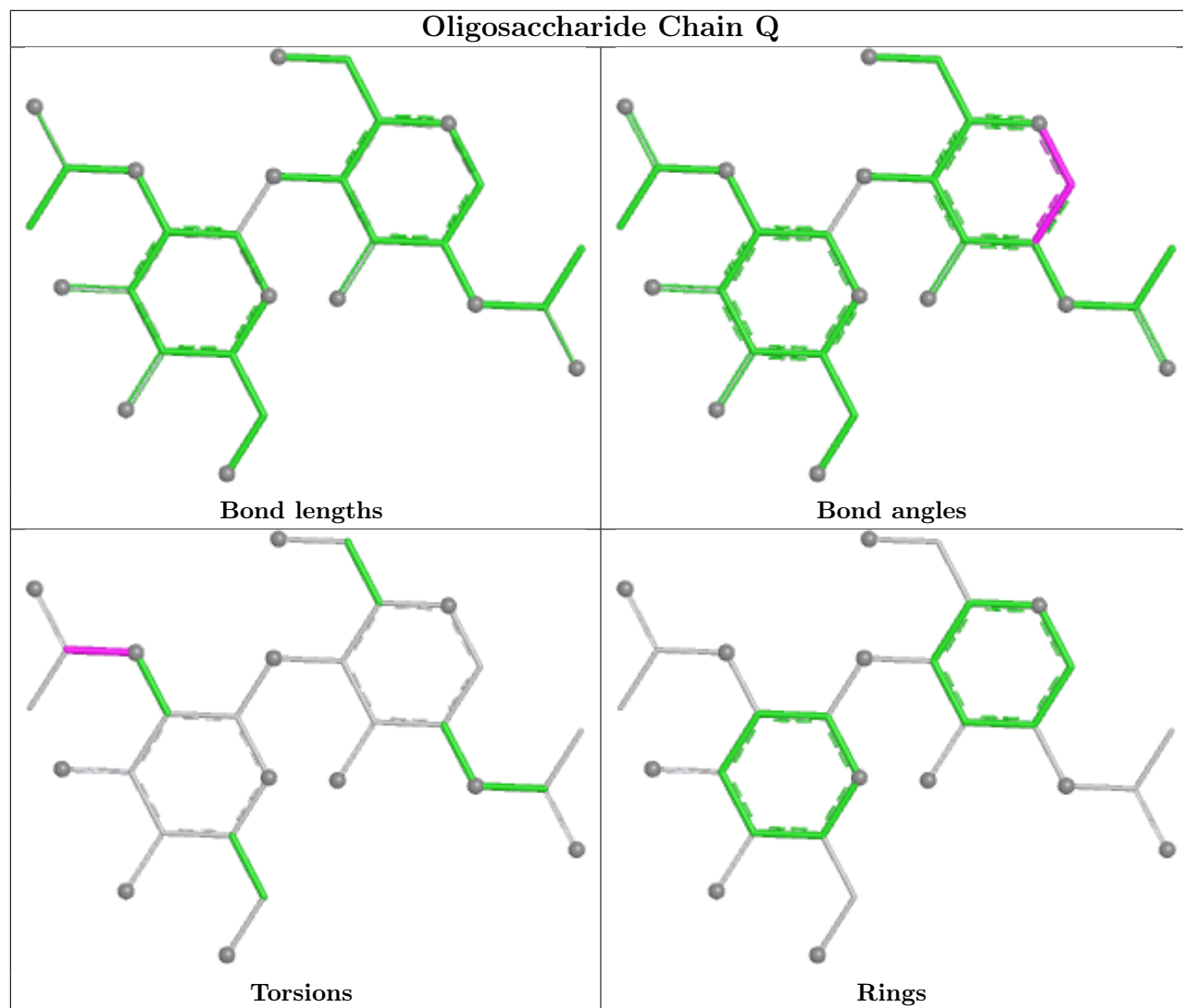


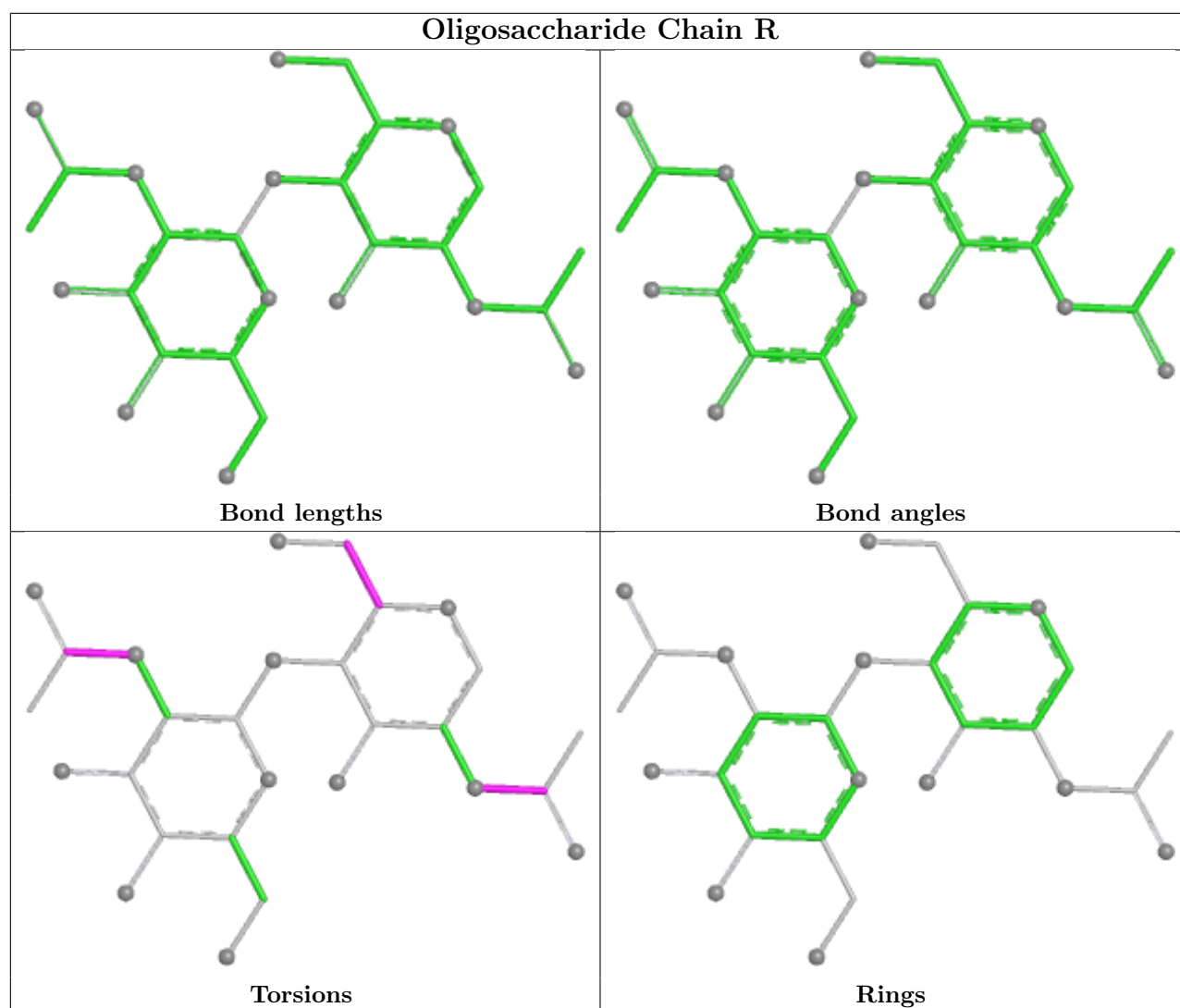


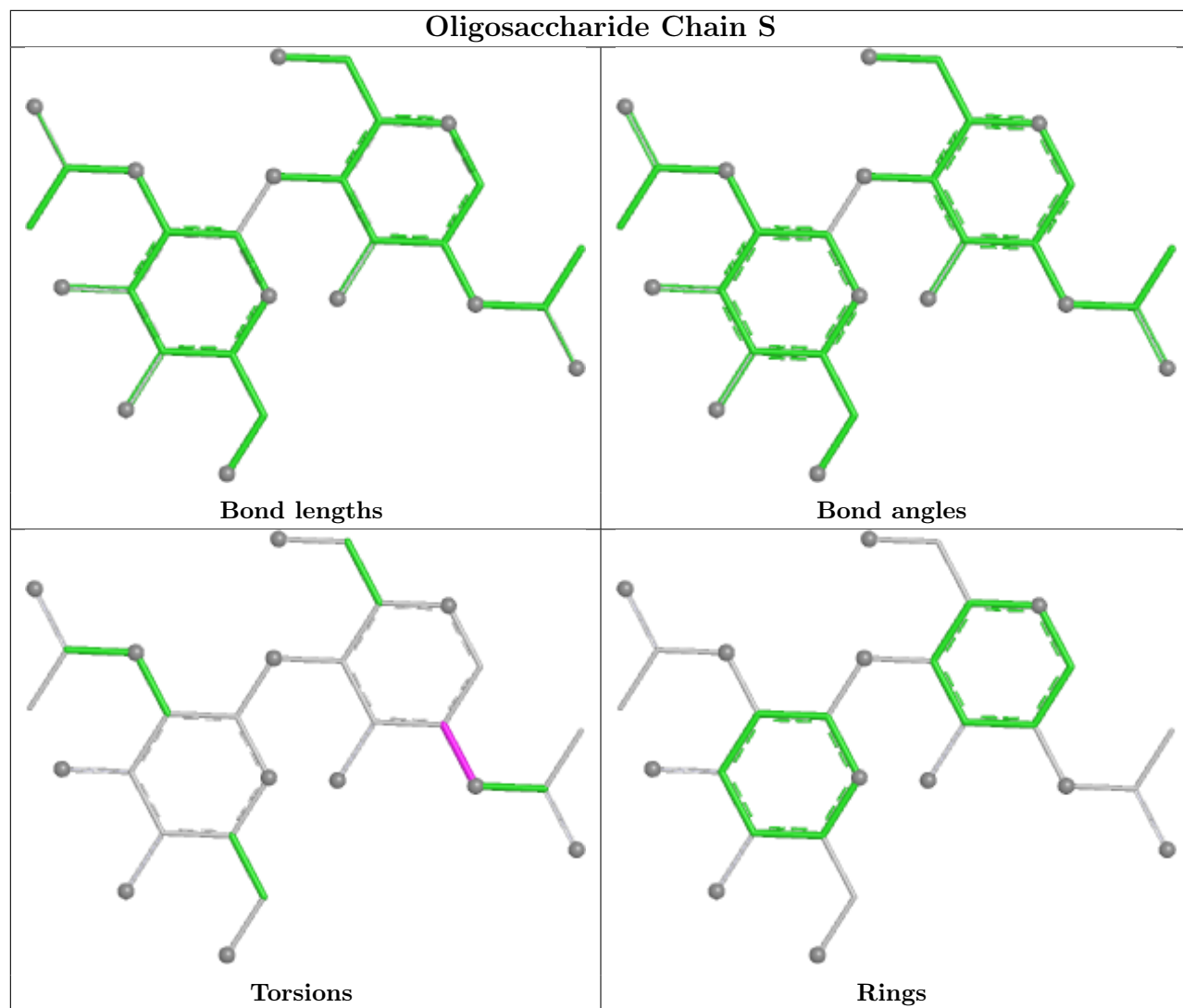


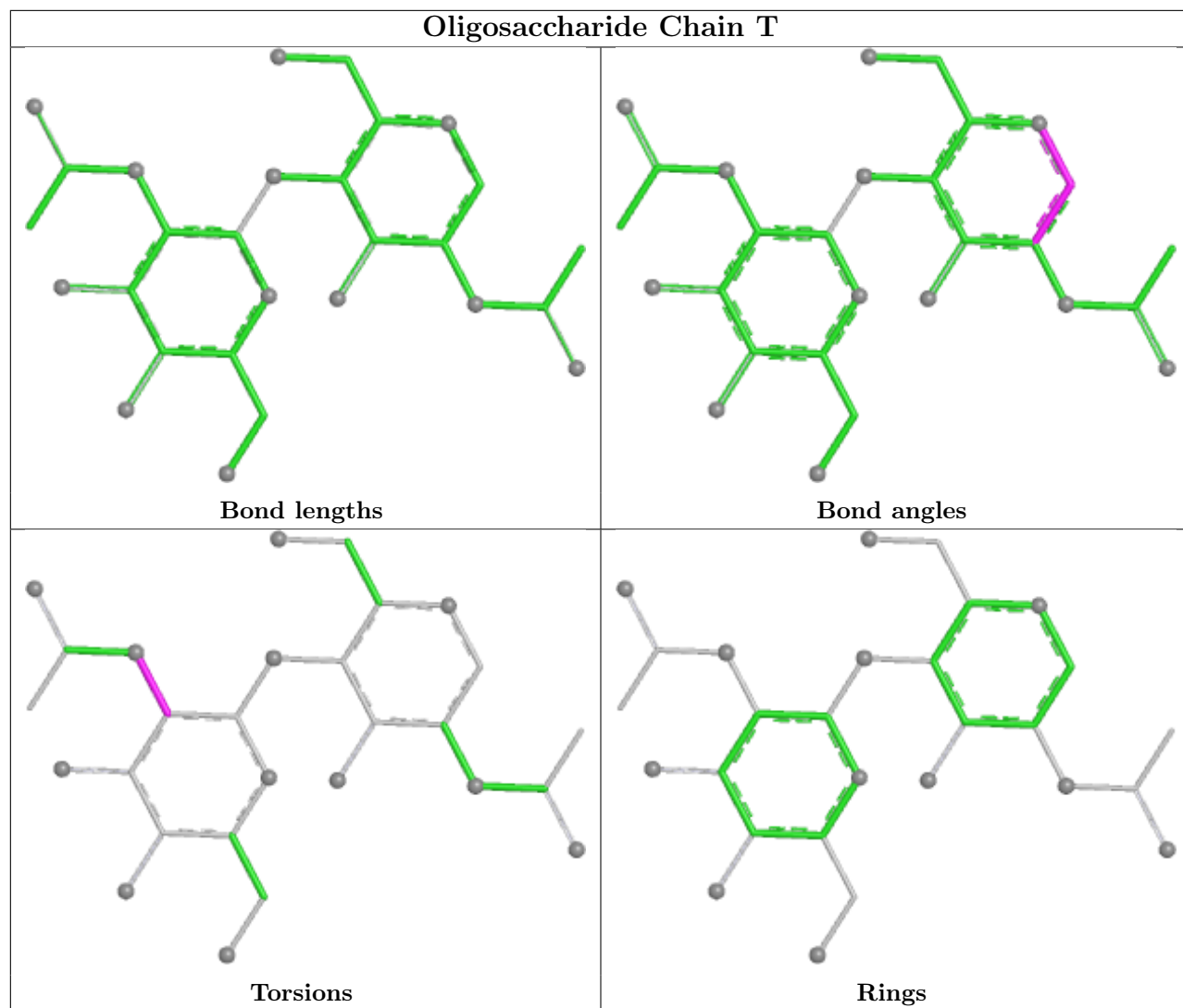


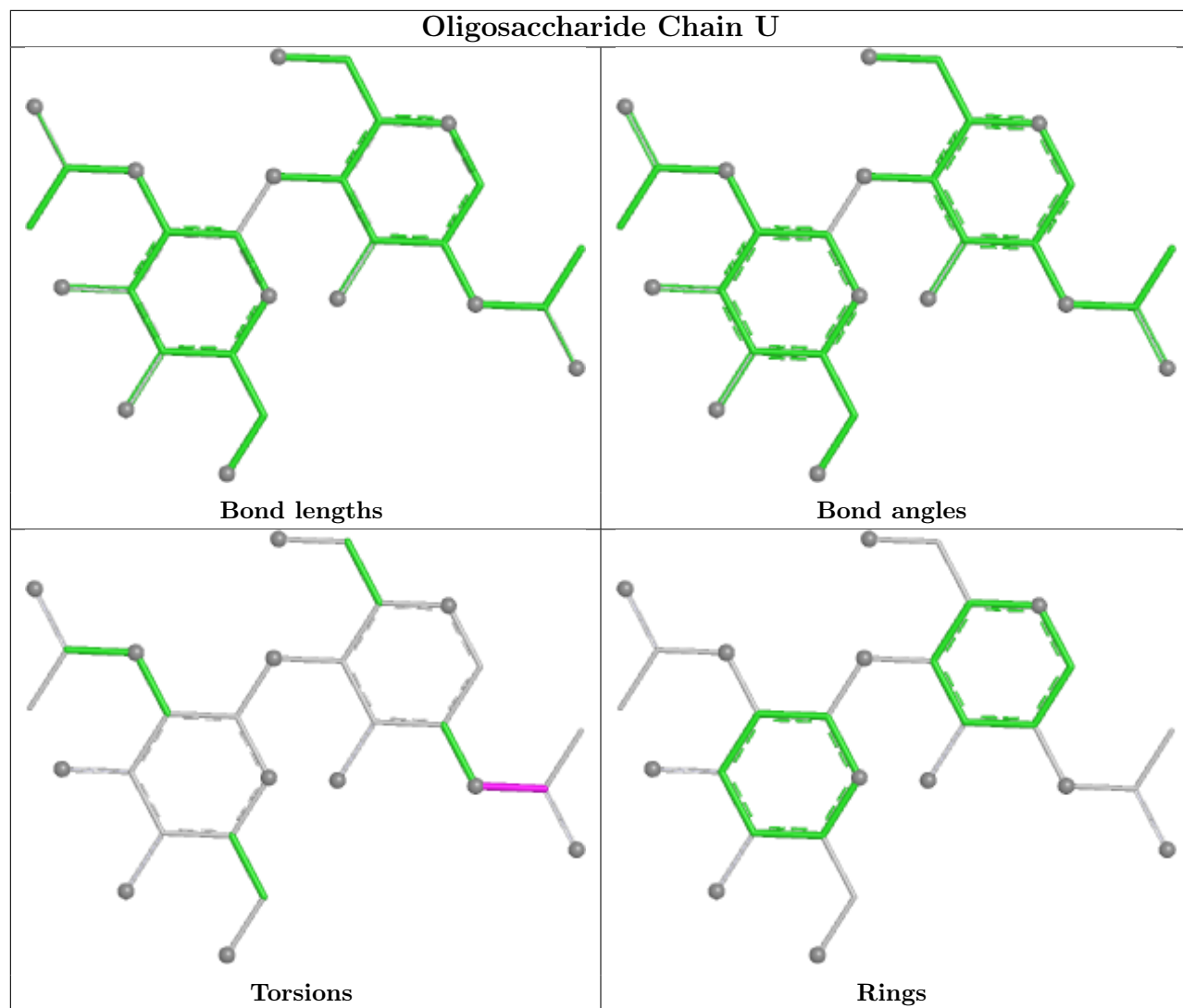


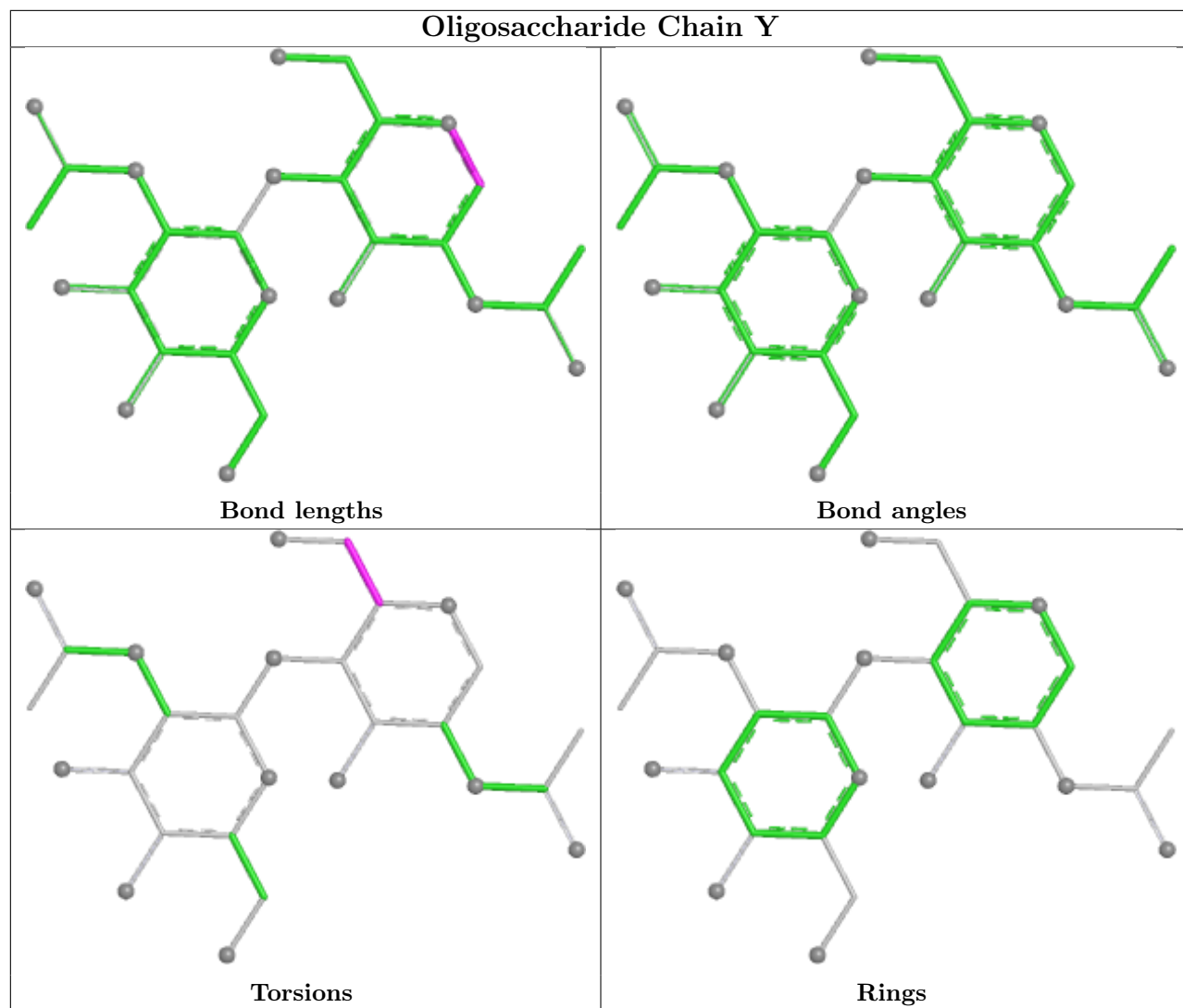


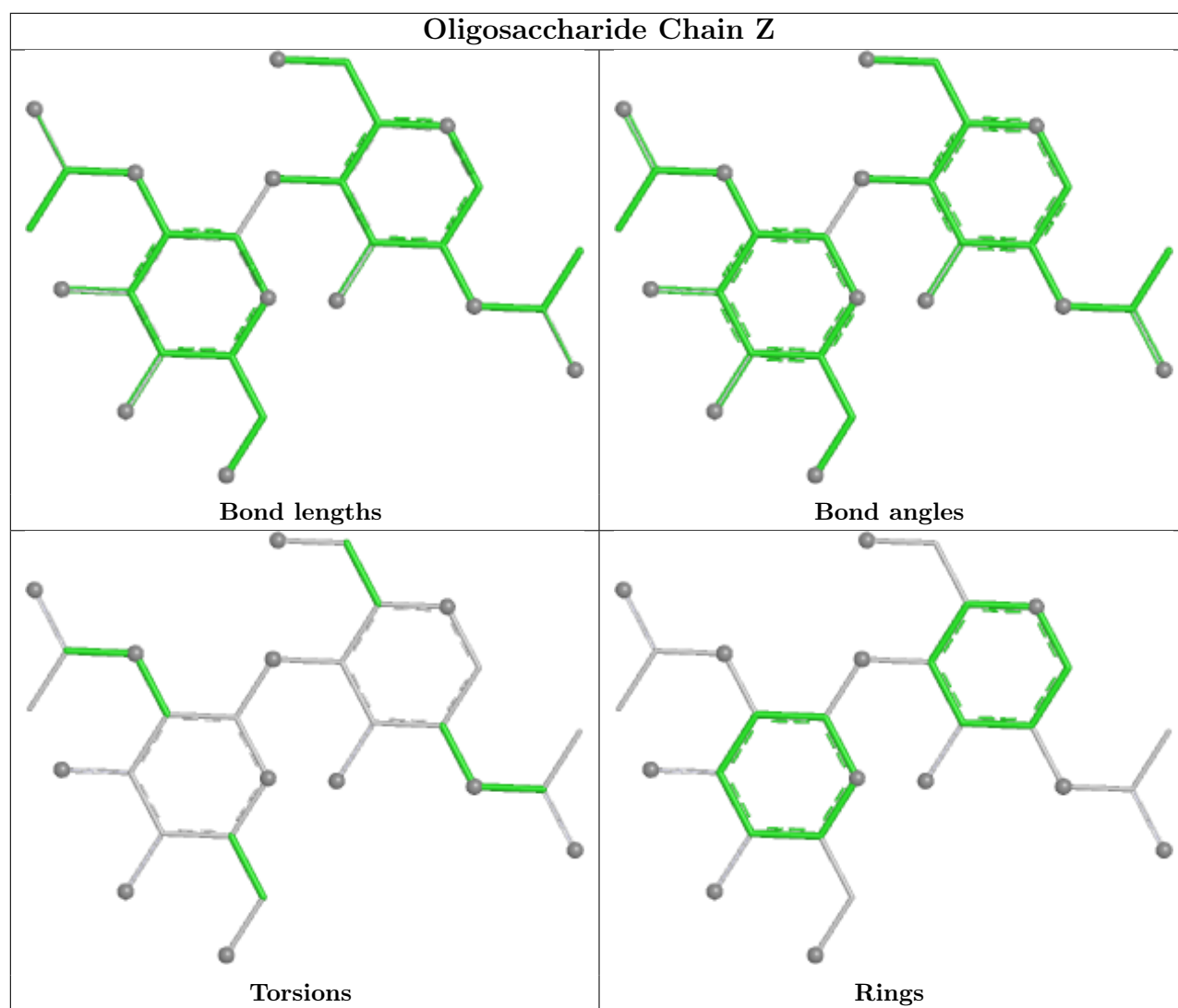


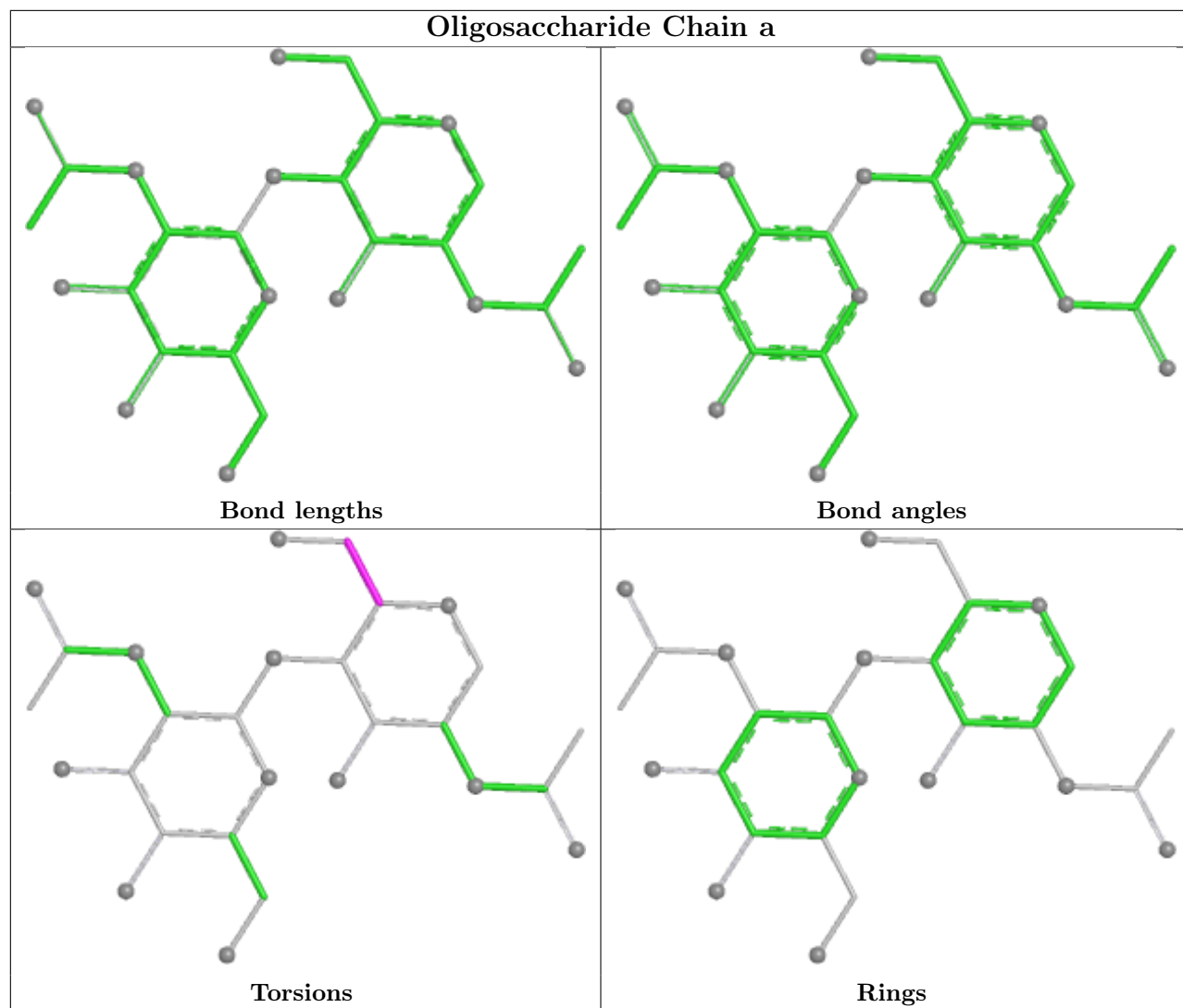


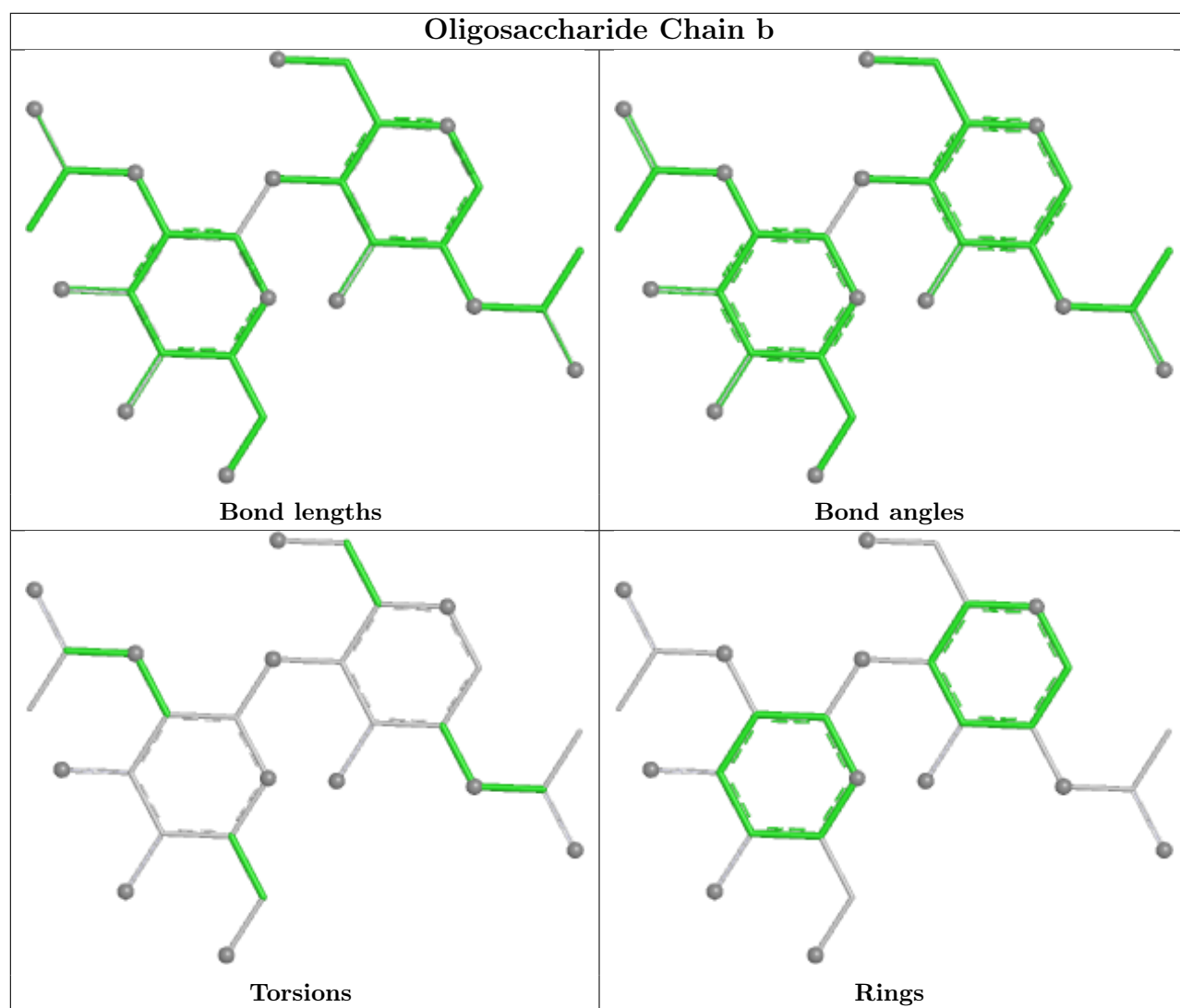


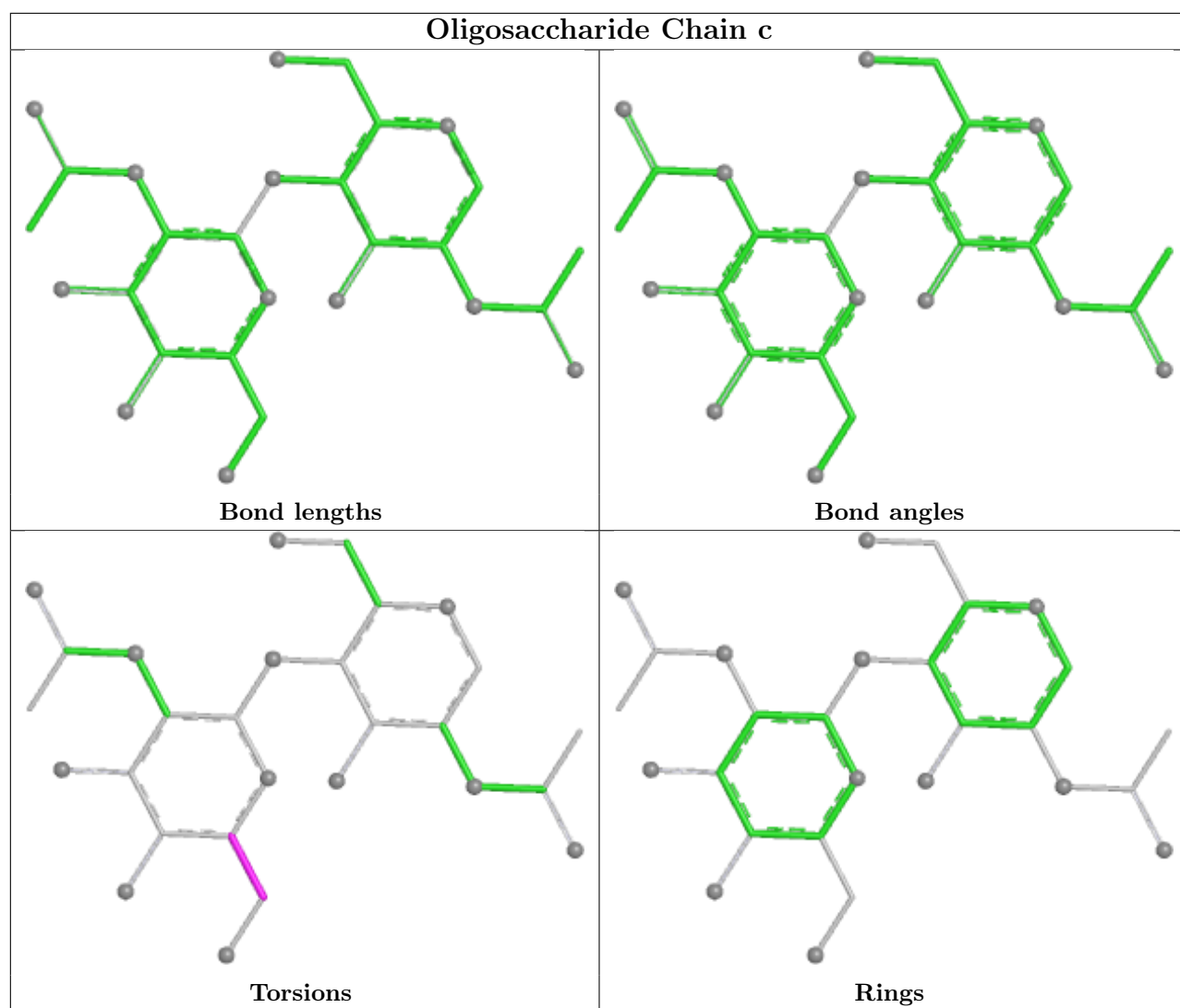


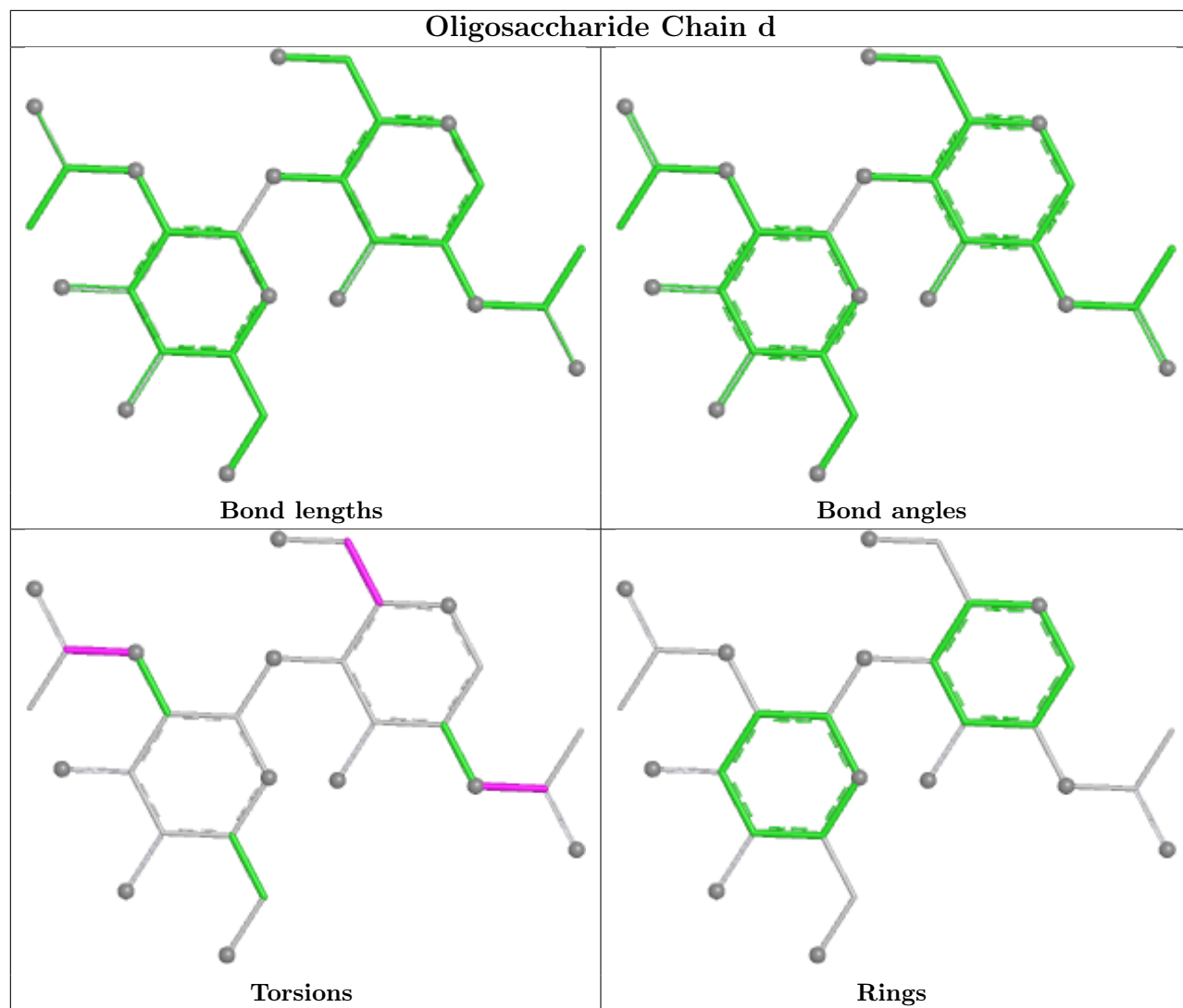


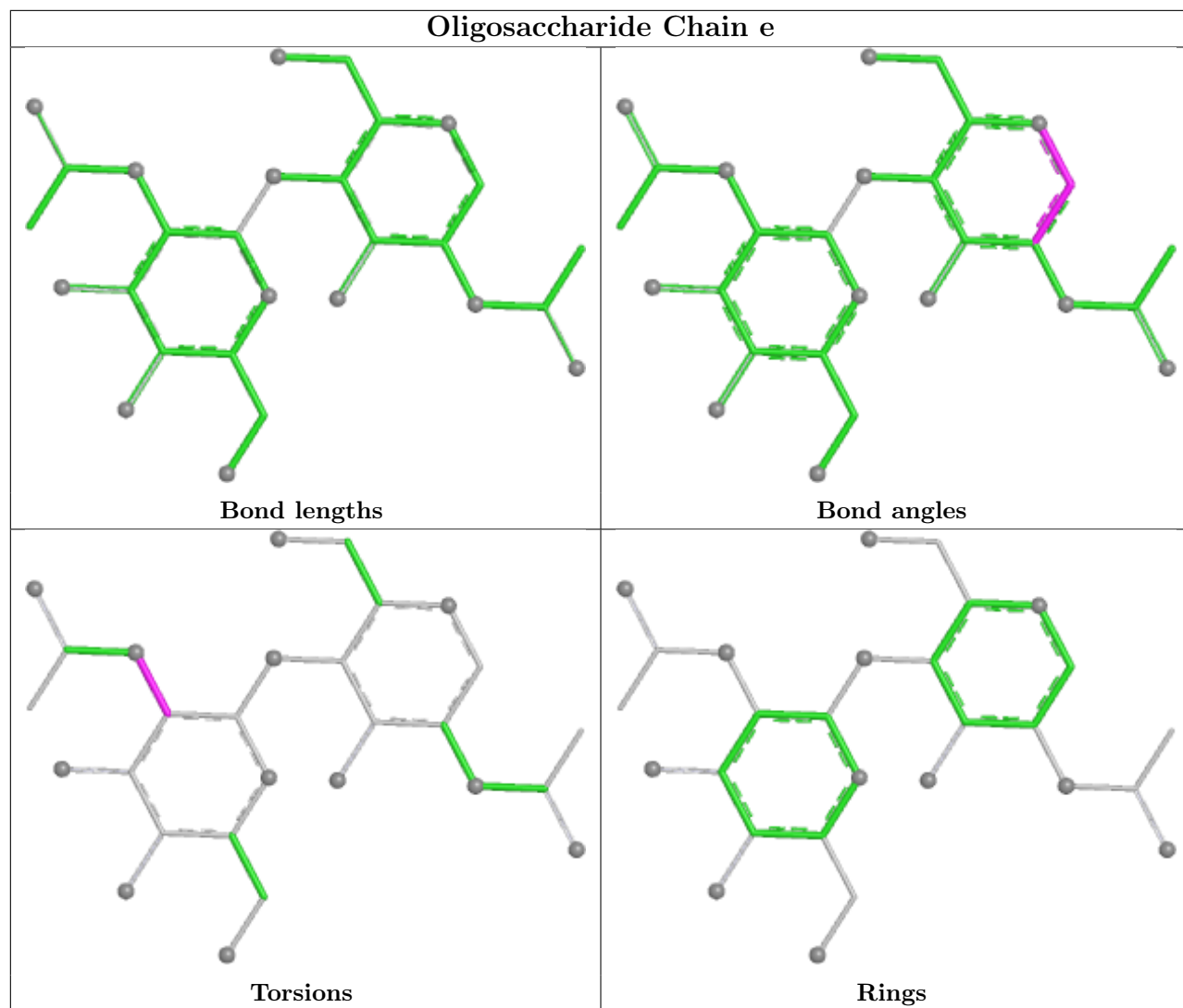


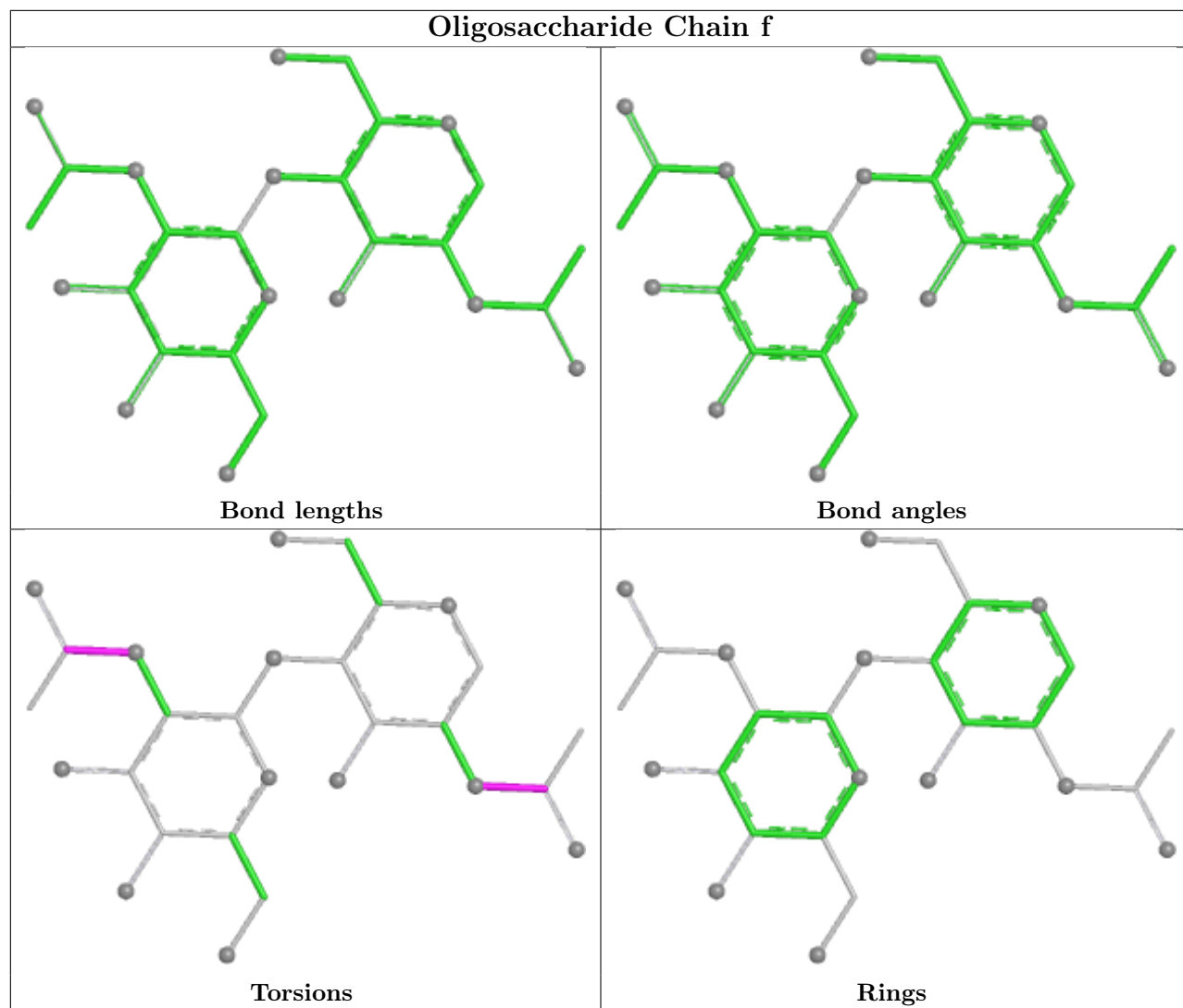


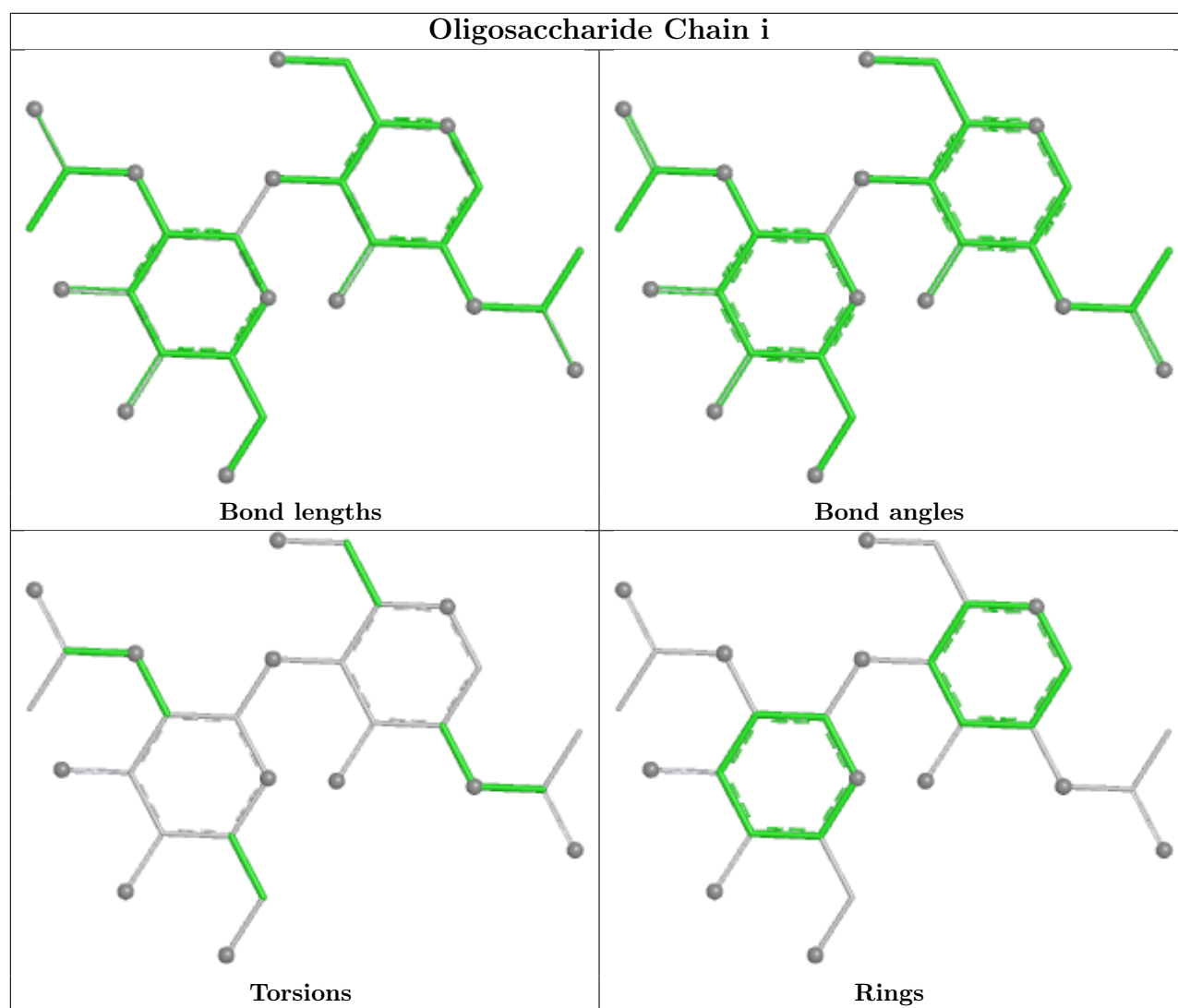


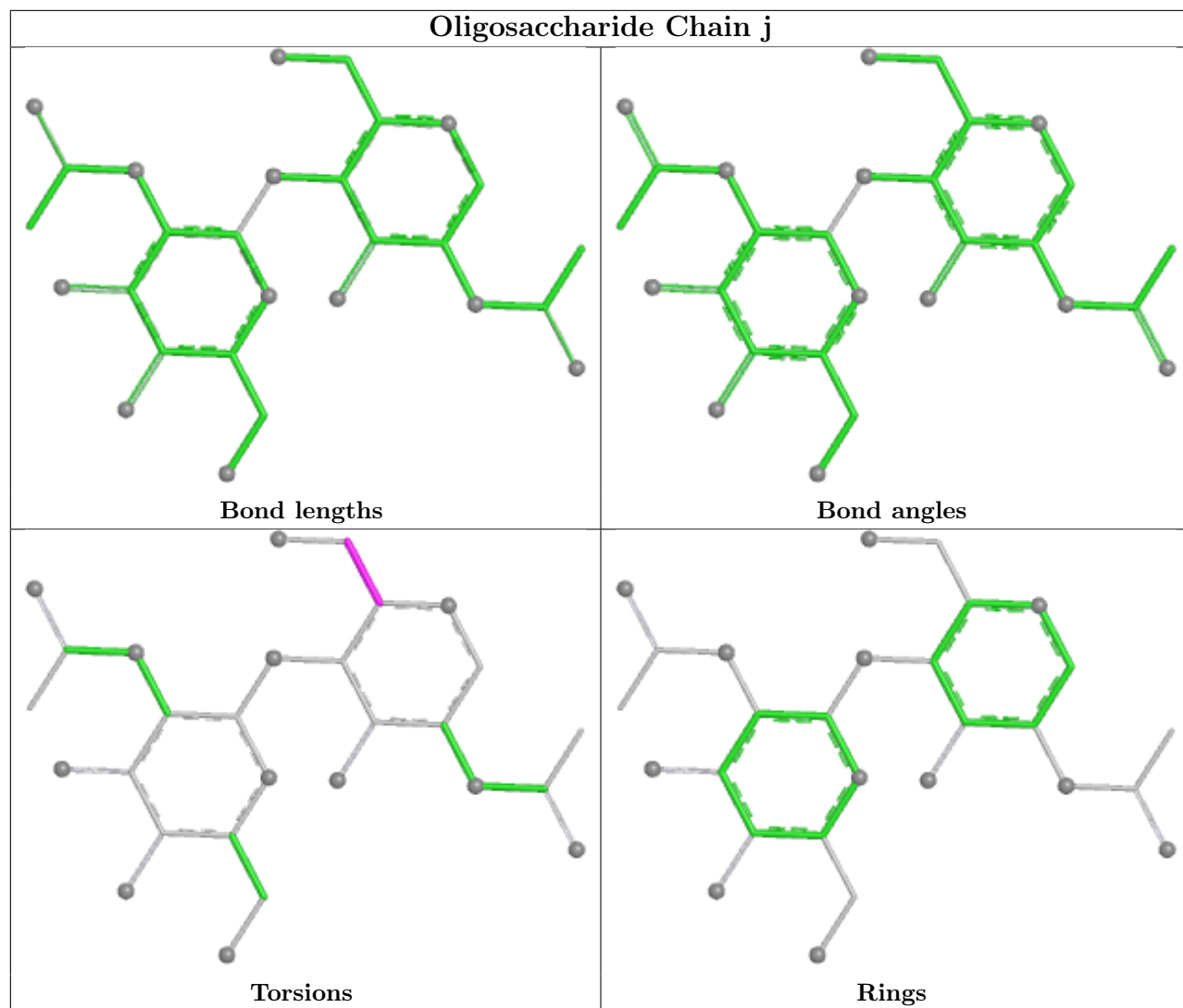


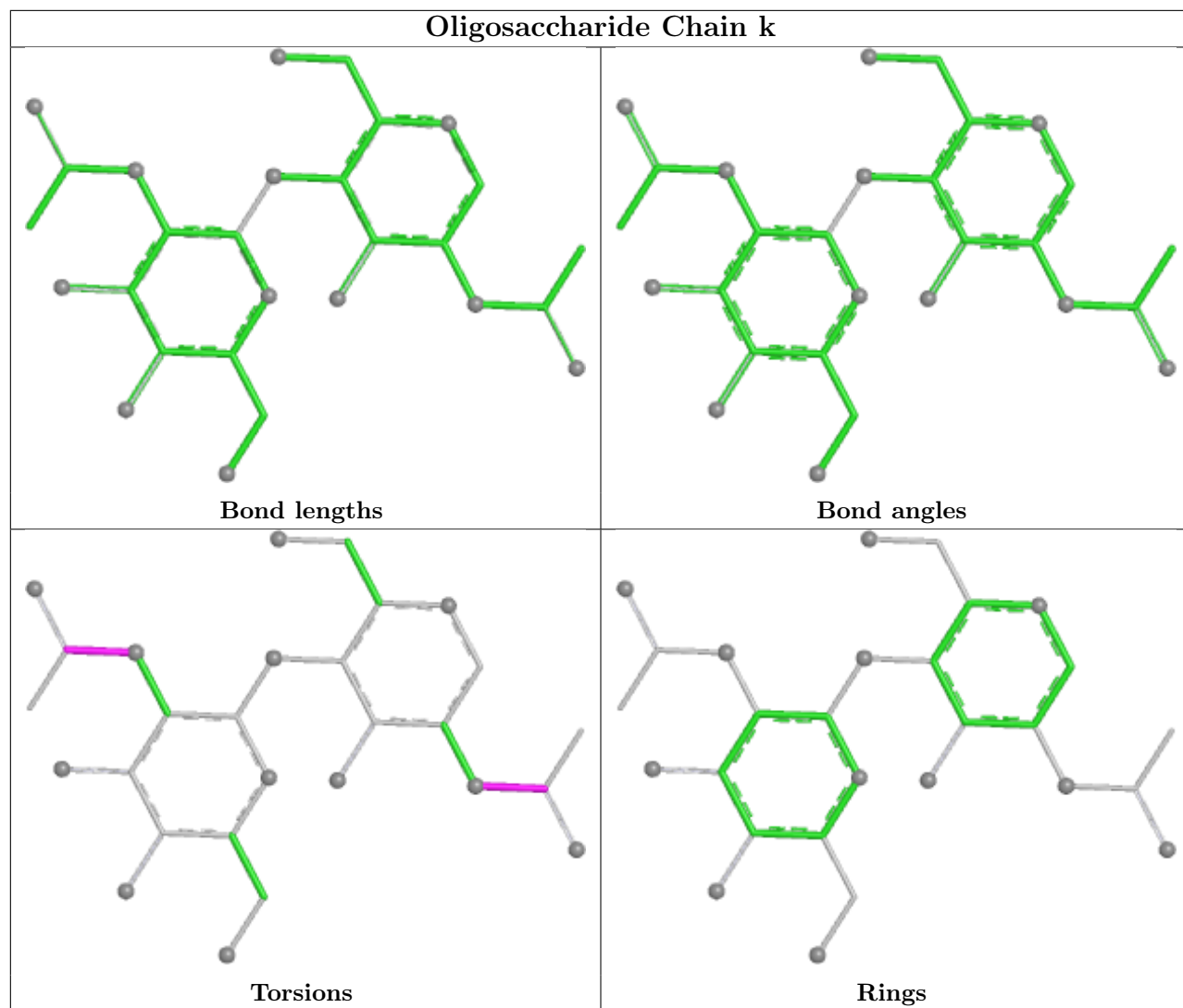


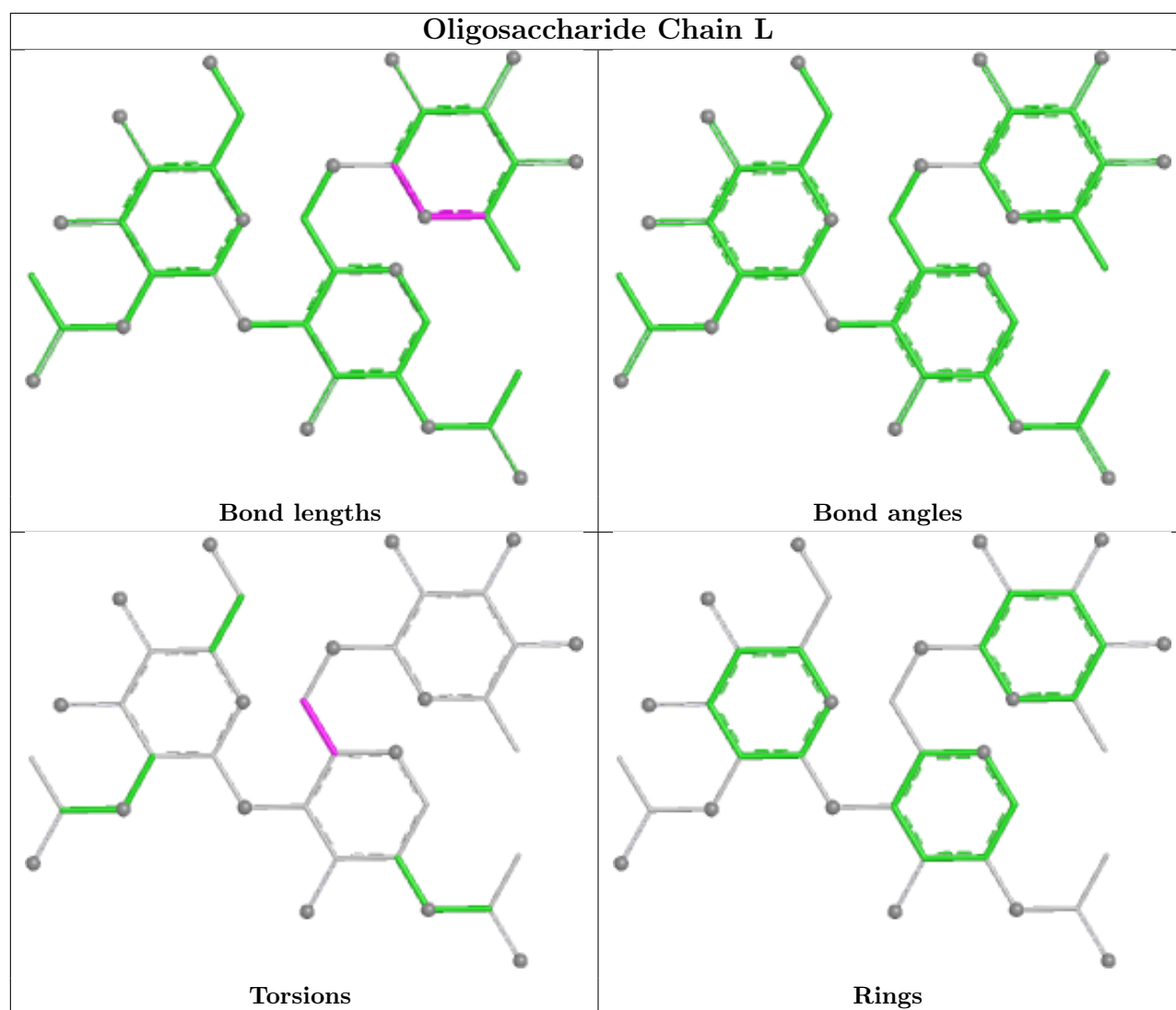


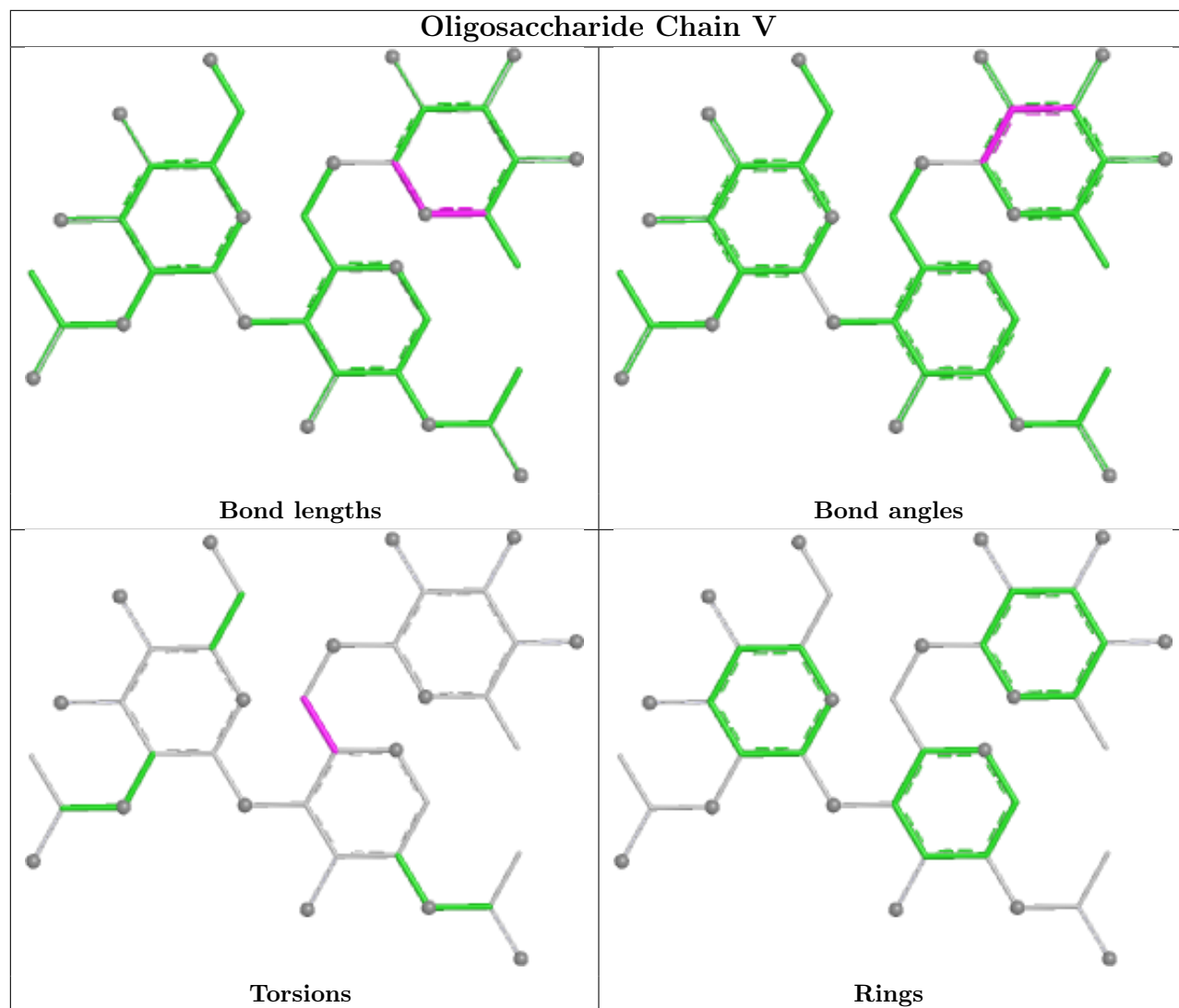


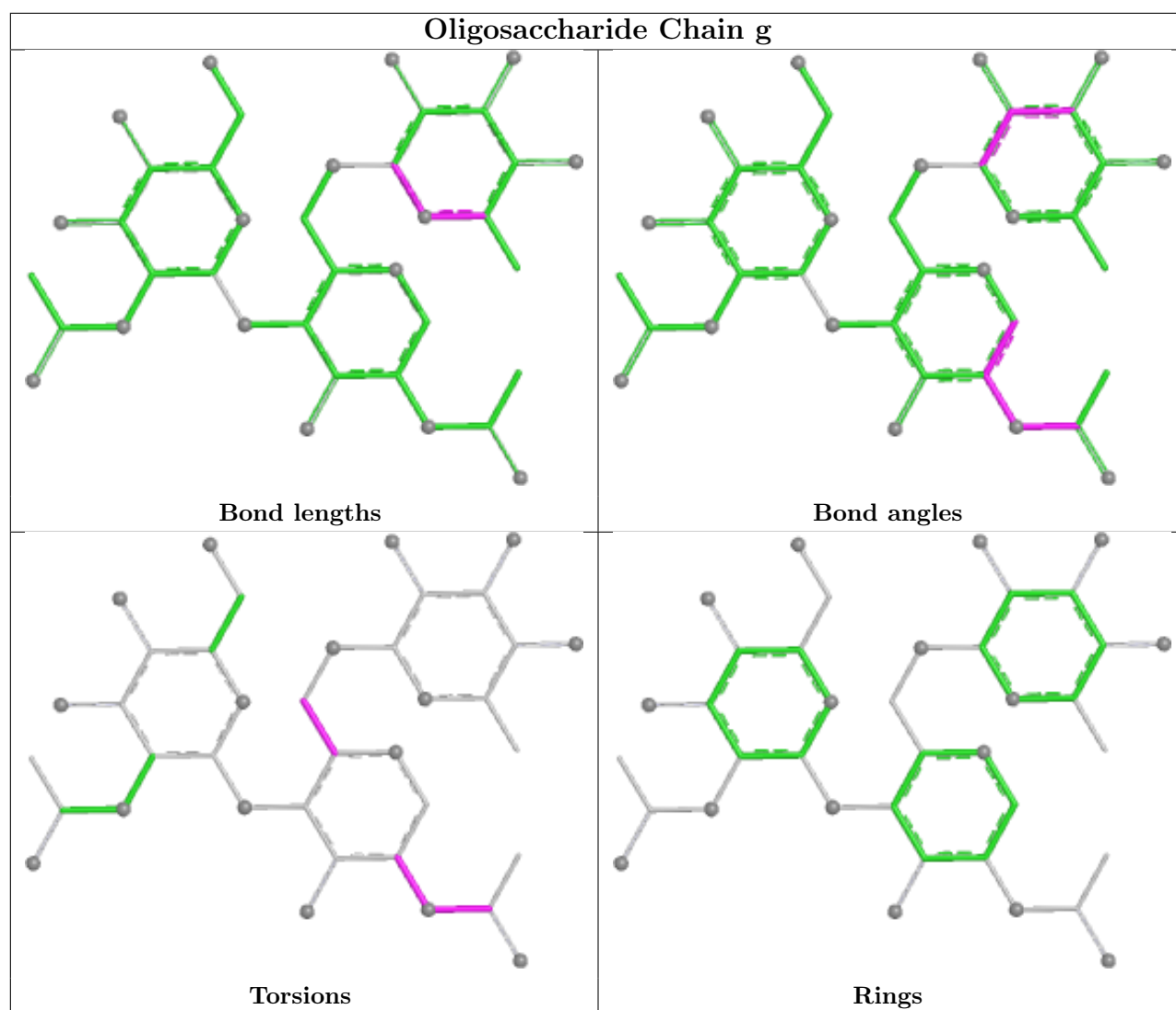


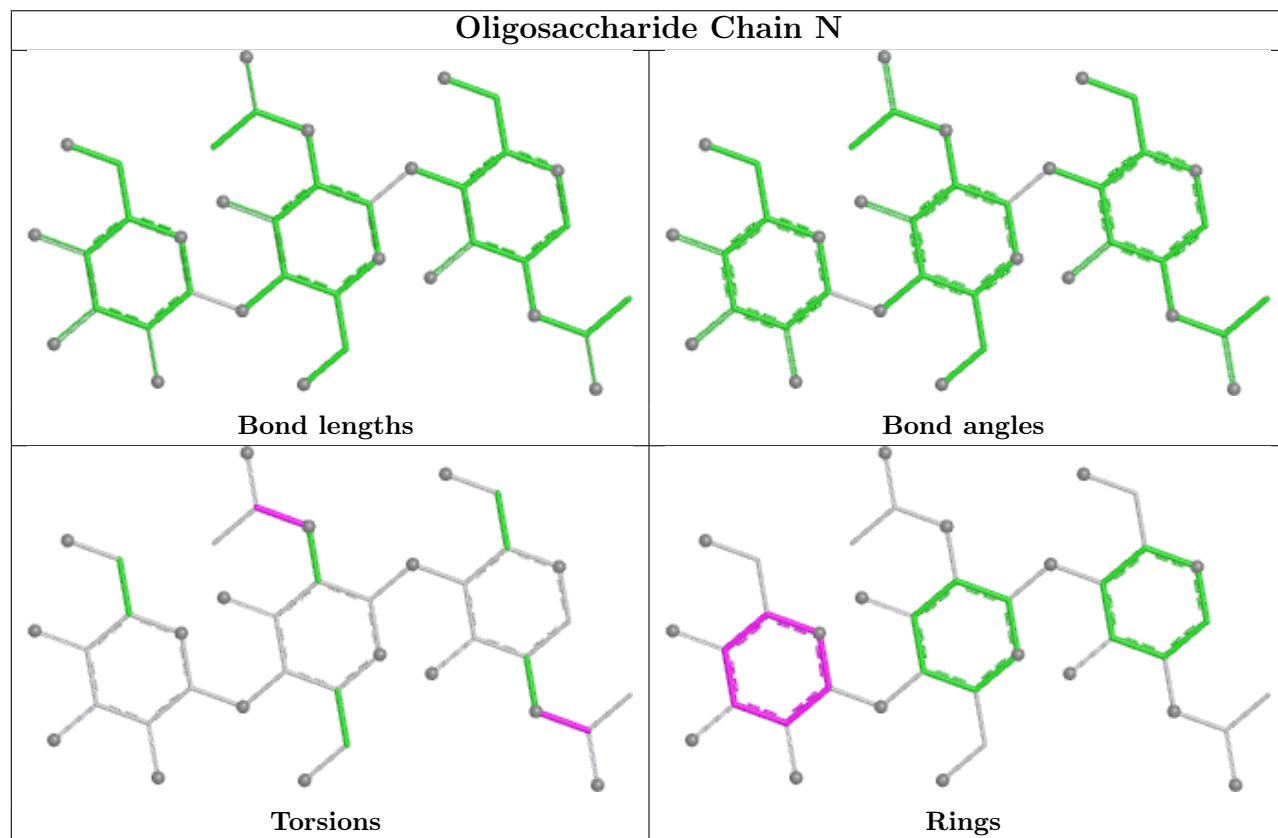
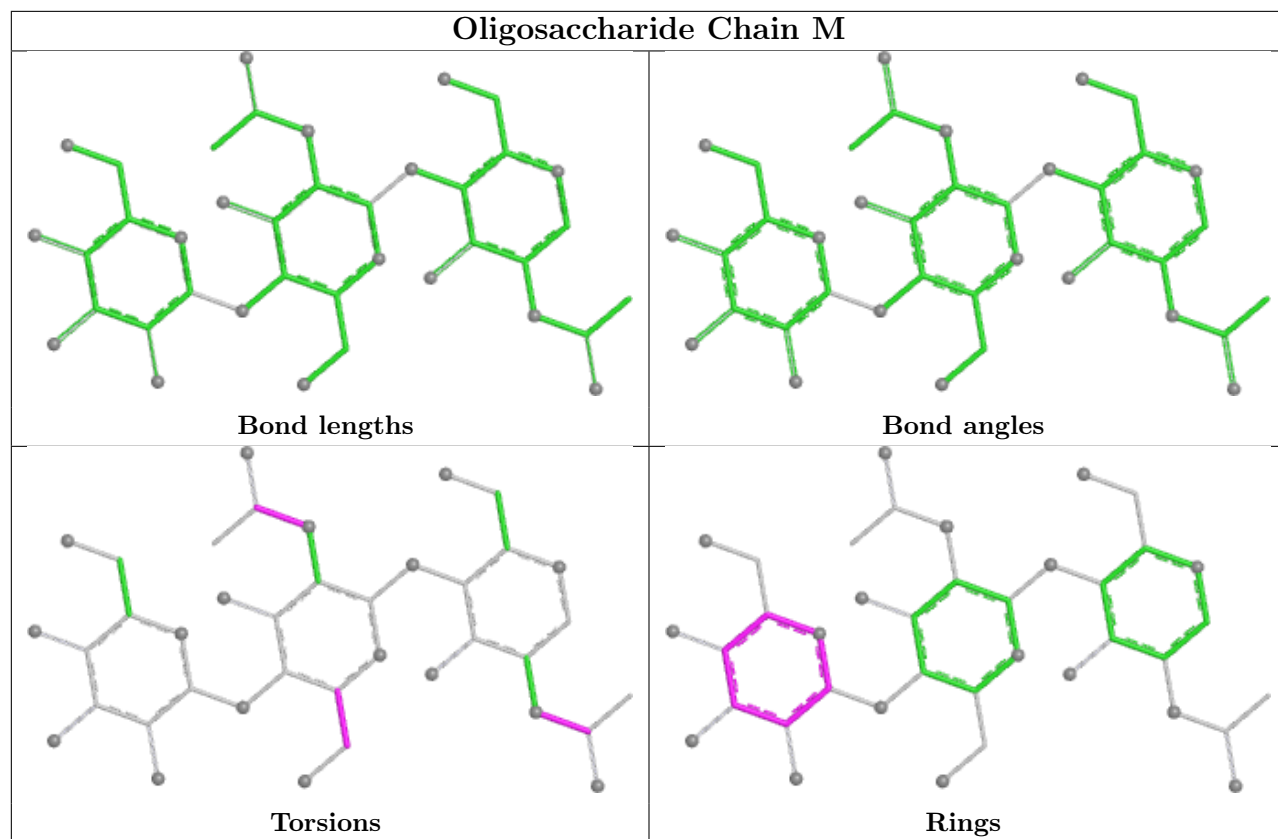


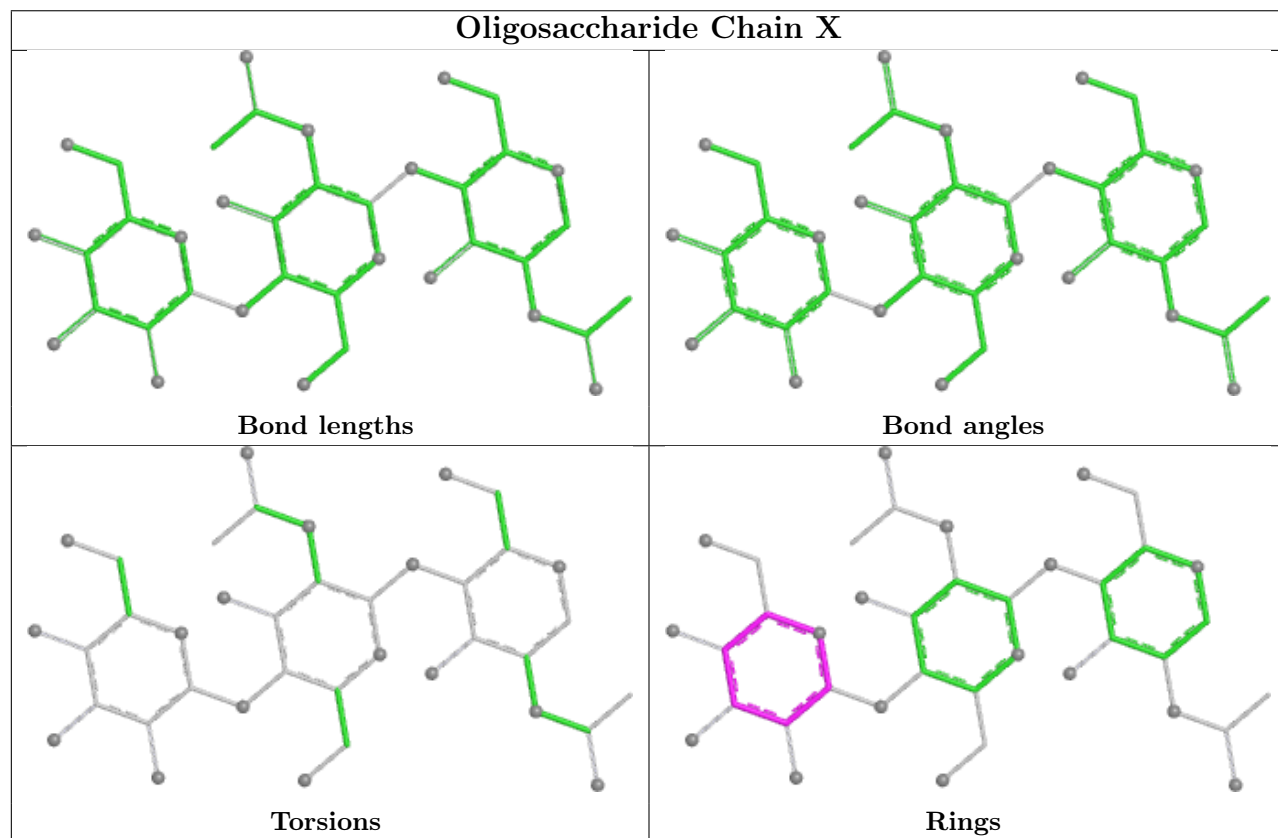
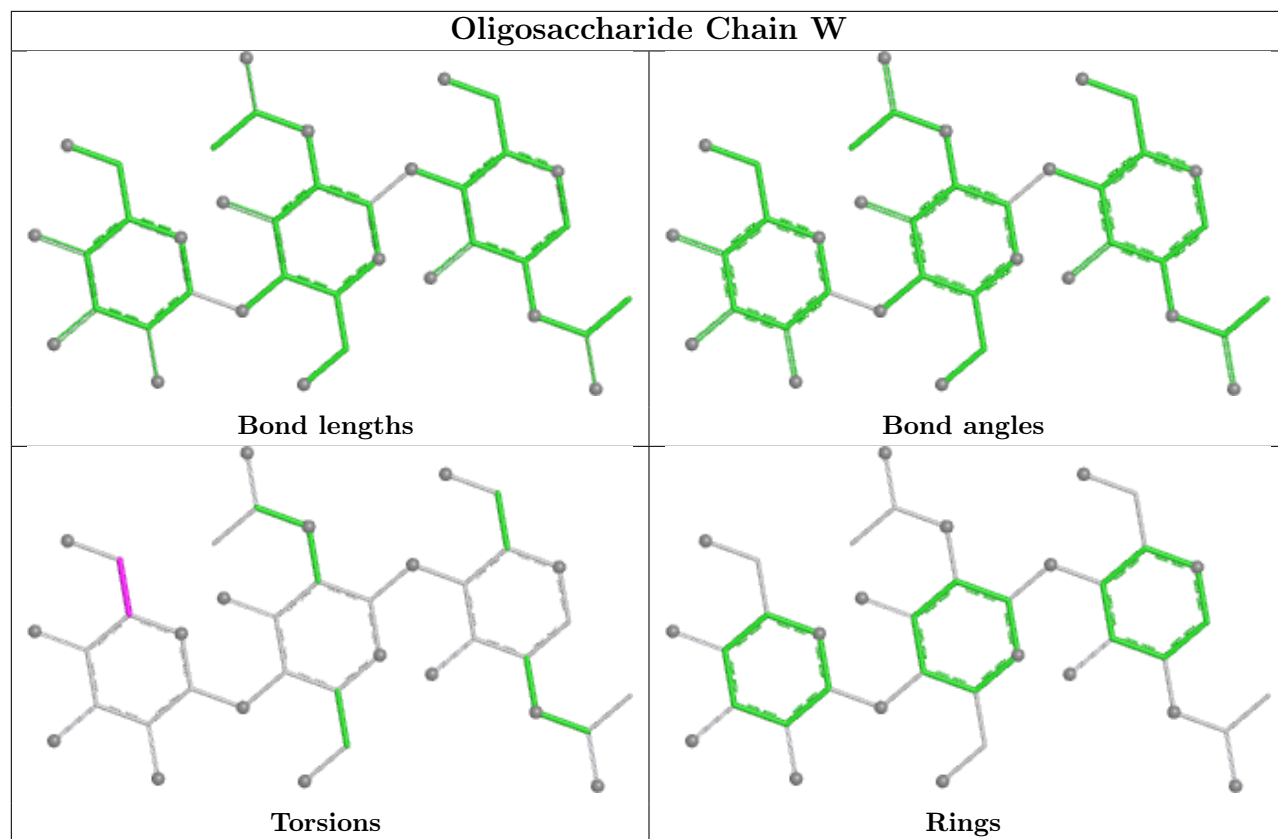


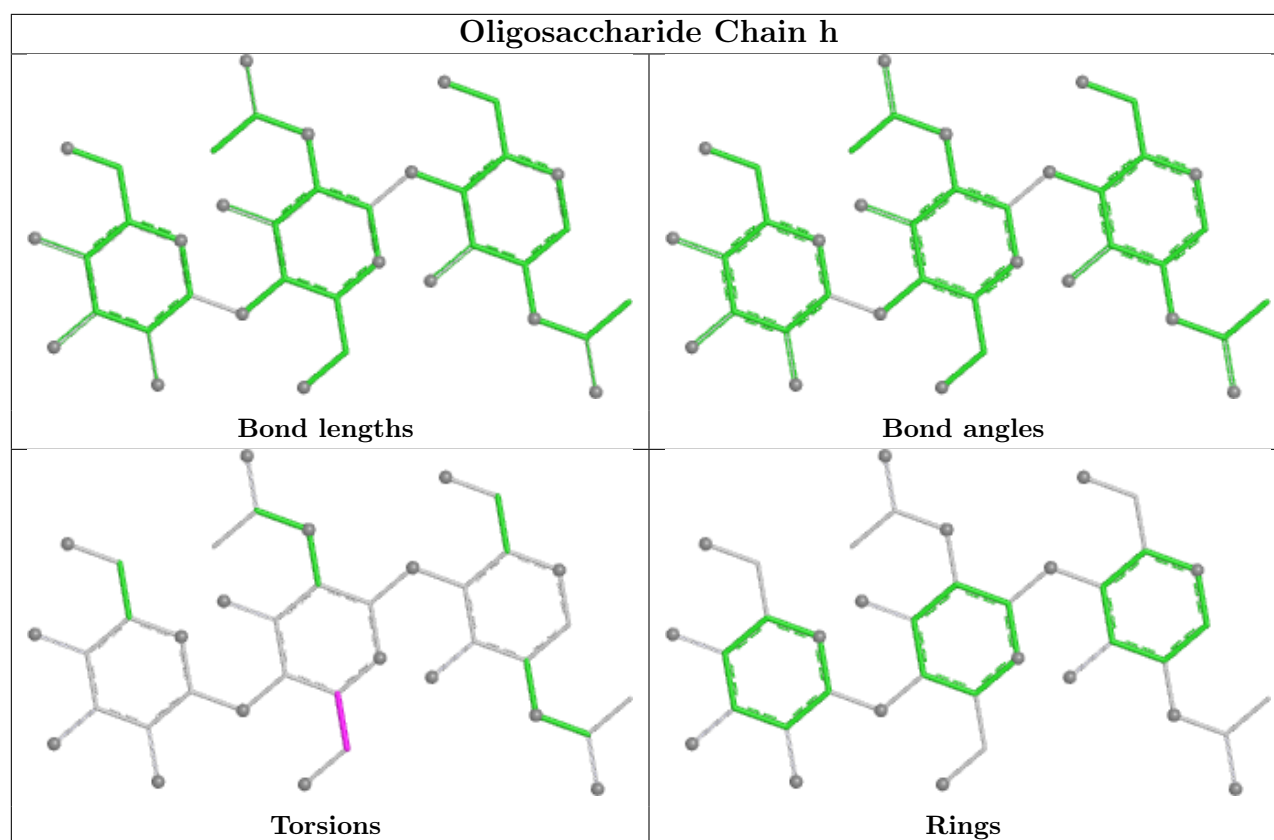












5.6 Ligand geometry [i](#)

25 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	$\# Z > 2$
7	NAG	A	1407	1	14,14,15	0.39	0	17,19,21	0.35	0
7	NAG	C	1407	1	14,14,15	0.40	0	17,19,21	0.34	0
7	NAG	A	1408	1	14,14,15	0.28	0	17,19,21	0.39	0
7	NAG	C	1403	1	14,14,15	0.40	0	17,19,21	0.57	0
7	NAG	B	1408	1	14,14,15	0.41	0	17,19,21	0.37	0
7	NAG	A	1409	1	14,14,15	0.39	0	17,19,21	0.37	0
7	NAG	B	1403	1	14,14,15	0.27	0	17,19,21	0.54	0
7	NAG	C	1408	1	14,14,15	0.17	0	17,19,21	0.46	0
7	NAG	A	1401	1	14,14,15	0.41	0	17,19,21	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NAG	C	1402	1	14,14,15	0.41	0	17,19,21	0.61	0
7	NAG	A	1403	1	14,14,15	0.41	0	17,19,21	0.33	0
7	NAG	C	1404	1	14,14,15	0.40	0	17,19,21	0.36	0
7	NAG	A	1404	1	14,14,15	0.25	0	17,19,21	0.39	0
7	NAG	B	1402	1	14,14,15	0.26	0	17,19,21	0.44	0
7	NAG	B	1407	1	14,14,15	0.40	0	17,19,21	0.38	0
7	NAG	B	1405	1	14,14,15	0.40	0	17,19,21	0.39	0
7	NAG	A	1406	1	14,14,15	0.41	0	17,19,21	0.38	0
7	NAG	A	1402	1	14,14,15	0.41	0	17,19,21	0.61	0
7	NAG	C	1406	1	14,14,15	0.41	0	17,19,21	0.41	0
7	NAG	B	1401	1	14,14,15	0.40	0	17,19,21	0.37	0
7	NAG	C	1401	1	14,14,15	0.42	0	17,19,21	0.47	0
7	NAG	B	1406	1	14,14,15	0.40	0	17,19,21	0.38	0
7	NAG	A	1405	1	14,14,15	0.40	0	17,19,21	0.34	0
7	NAG	B	1404	1	14,14,15	0.75	1 (7%)	17,19,21	0.71	1 (5%)
7	NAG	C	1405	1	14,14,15	0.38	0	17,19,21	0.35	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
7	NAG	A	1407	1	-	2/6/23/26	0/1/1/1
7	NAG	C	1407	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1408	1	-	0/6/23/26	0/1/1/1
7	NAG	C	1403	1	-	4/6/23/26	0/1/1/1
7	NAG	B	1408	1	-	3/6/23/26	0/1/1/1
7	NAG	A	1409	1	-	0/6/23/26	0/1/1/1
7	NAG	B	1403	1	-	2/6/23/26	0/1/1/1
7	NAG	C	1408	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1401	1	-	1/6/23/26	0/1/1/1
7	NAG	C	1402	1	-	4/6/23/26	0/1/1/1
7	NAG	A	1403	1	-	0/6/23/26	0/1/1/1
7	NAG	C	1404	1	-	0/6/23/26	0/1/1/1
7	NAG	A	1404	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1402	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1407	1	-	0/6/23/26	0/1/1/1
7	NAG	B	1405	1	-	3/6/23/26	0/1/1/1
7	NAG	A	1406	1	-	3/6/23/26	0/1/1/1

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	1404	NAG	O5-C1	-2.65	1.39	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	1404	NAG	C2-N2-C7	2.23	125.89	122.90

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1402	NAG	C8-C7-N2-C2
7	A	1402	NAG	O7-C7-N2-C2
7	A	1405	NAG	C8-C7-N2-C2
7	A	1405	NAG	O7-C7-N2-C2
7	C	1402	NAG	C8-C7-N2-C2
7	C	1402	NAG	O7-C7-N2-C2
7	C	1403	NAG	C8-C7-N2-C2
7	C	1403	NAG	O7-C7-N2-C2
7	C	1405	NAG	C8-C7-N2-C2
7	C	1405	NAG	O7-C7-N2-C2
7	C	1407	NAG	C8-C7-N2-C2
7	C	1407	NAG	O7-C7-N2-C2
7	B	1402	NAG	O5-C5-C6-O6
7	A	1407	NAG	C8-C7-N2-C2
7	A	1407	NAG	O7-C7-N2-C2
7	C	1408	NAG	O5-C5-C6-O6
7	A	1404	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	B	1402	NAG	C4-C5-C6-O6
7	B	1405	NAG	C8-C7-N2-C2
7	C	1405	NAG	O5-C5-C6-O6
7	C	1408	NAG	C4-C5-C6-O6
7	A	1404	NAG	C4-C5-C6-O6
7	C	1405	NAG	C4-C5-C6-O6
7	A	1402	NAG	O5-C5-C6-O6
7	C	1402	NAG	O5-C5-C6-O6
7	B	1401	NAG	C8-C7-N2-C2
7	B	1405	NAG	O7-C7-N2-C2
7	B	1403	NAG	O5-C5-C6-O6
7	B	1401	NAG	O7-C7-N2-C2
7	B	1403	NAG	C4-C5-C6-O6
7	A	1402	NAG	C4-C5-C6-O6
7	C	1402	NAG	C4-C5-C6-O6
7	A	1406	NAG	C8-C7-N2-C2
7	B	1408	NAG	C8-C7-N2-C2
7	C	1403	NAG	O5-C5-C6-O6
7	A	1401	NAG	O5-C5-C6-O6
7	B	1404	NAG	O5-C5-C6-O6
7	A	1406	NAG	O7-C7-N2-C2
7	B	1408	NAG	O7-C7-N2-C2
7	A	1406	NAG	C1-C2-N2-C7
7	B	1405	NAG	C1-C2-N2-C7
7	B	1408	NAG	C1-C2-N2-C7
7	C	1403	NAG	C3-C2-N2-C7

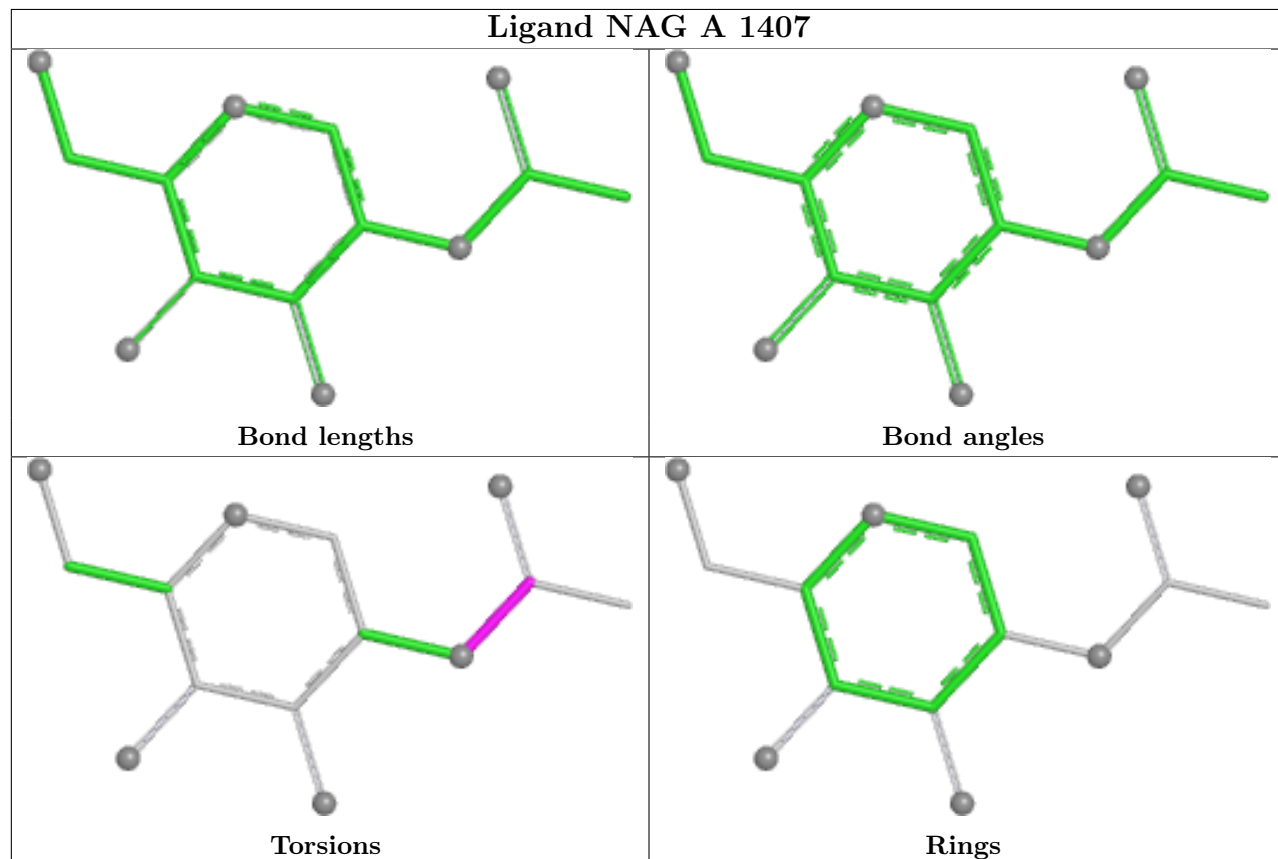
There are no ring outliers.

3 monomers are involved in 3 short contacts:

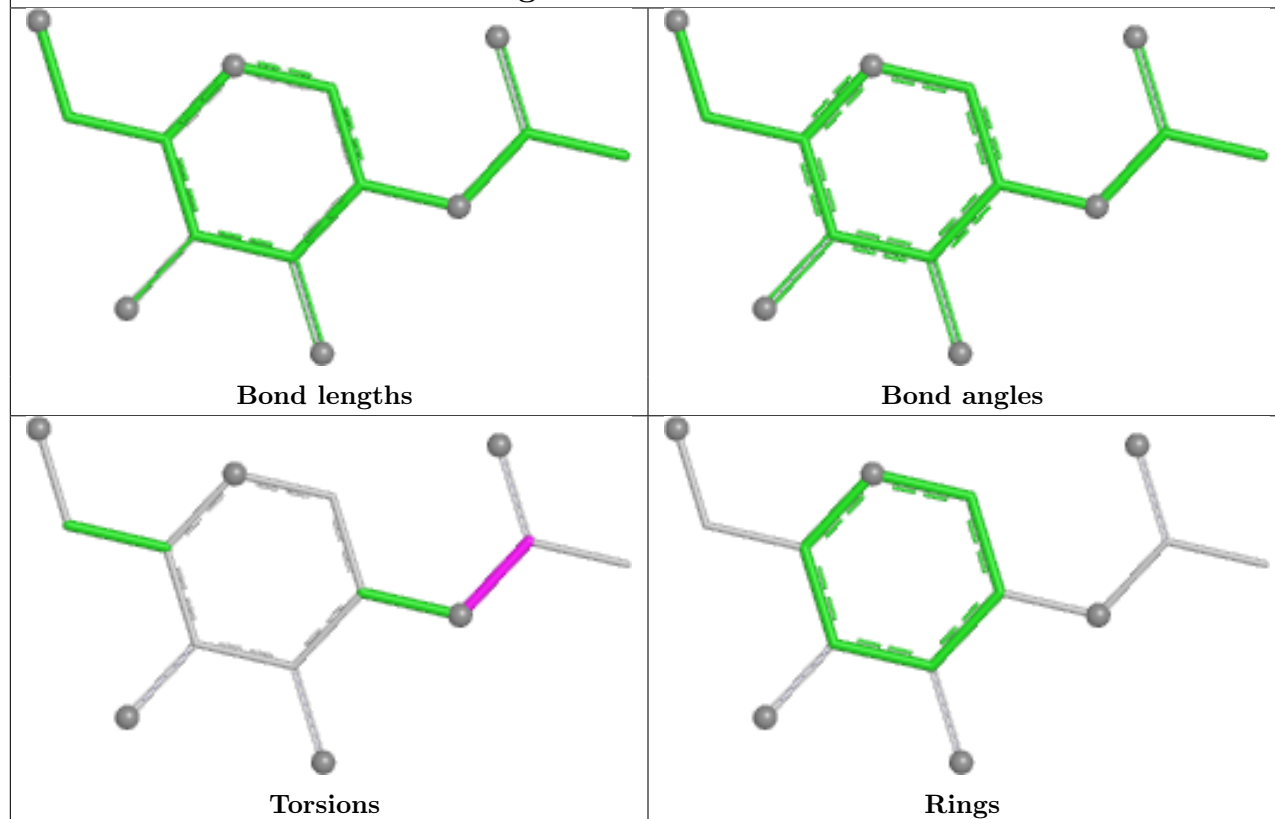
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	1402	NAG	1	0
7	A	1402	NAG	1	0
7	C	1401	NAG	1	0

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

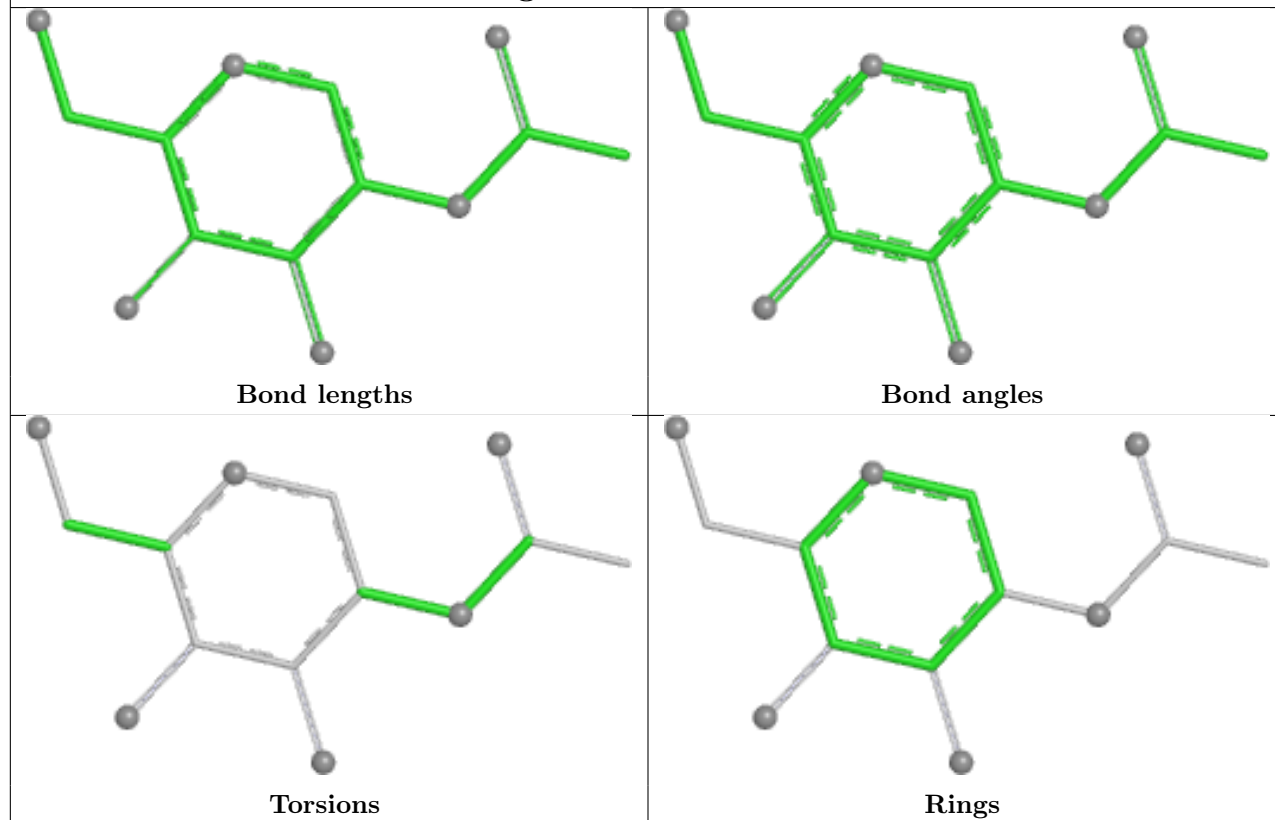
in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



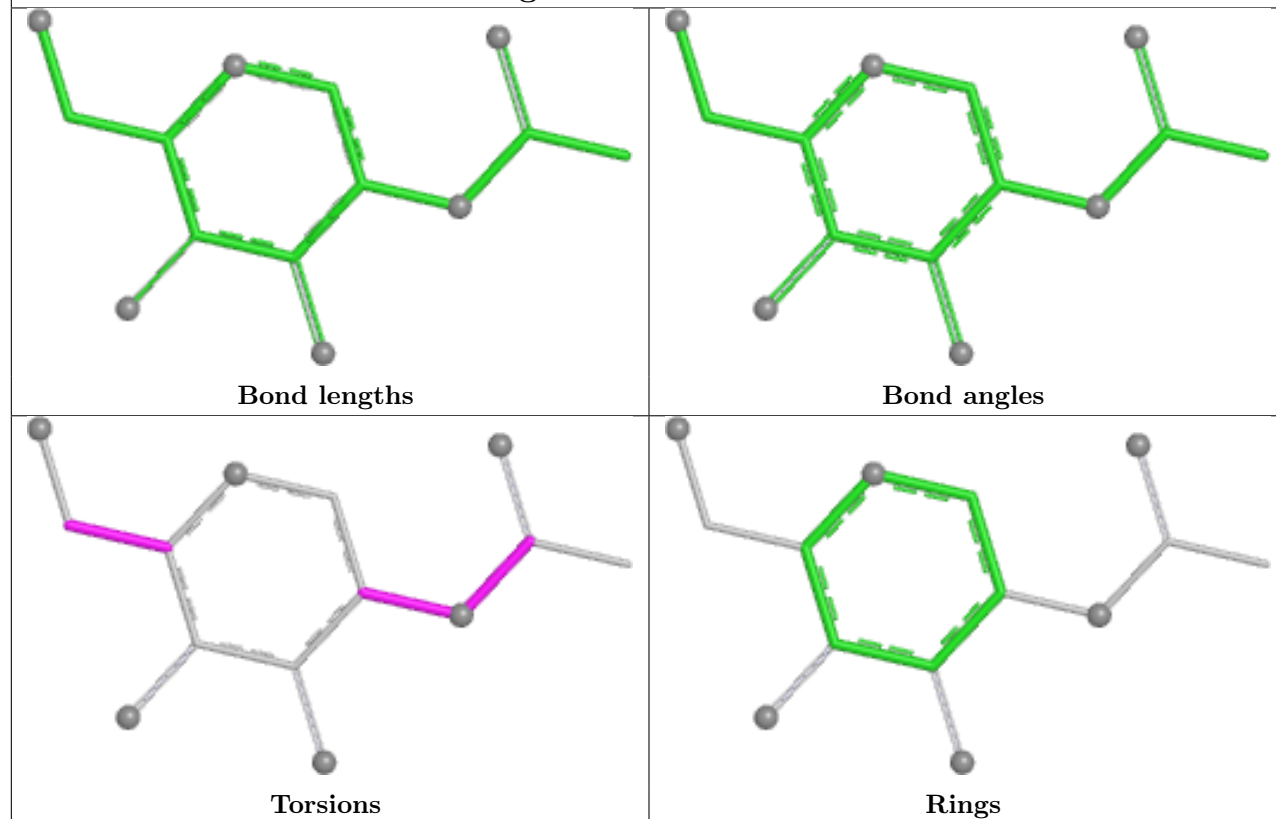
Ligand NAG C 1407



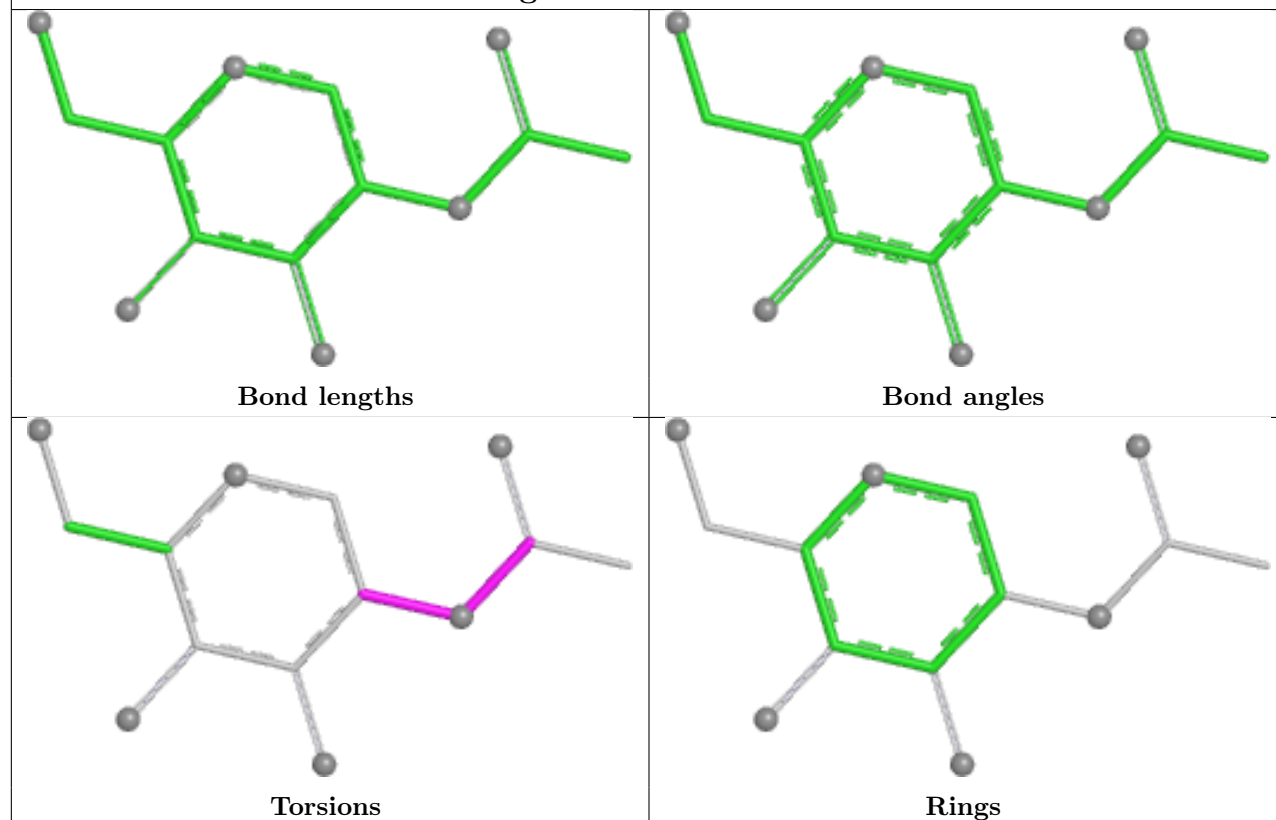
Ligand NAG A 1408



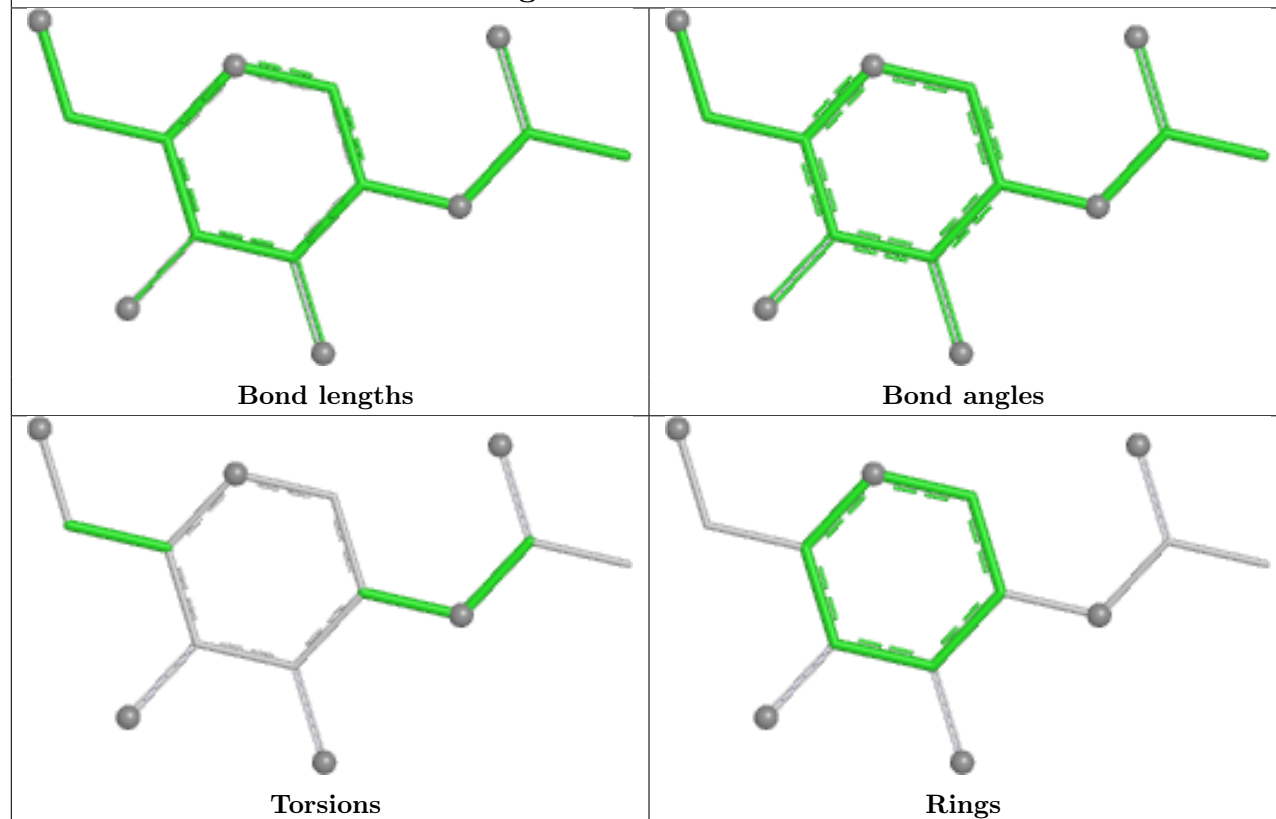
Ligand NAG C 1403



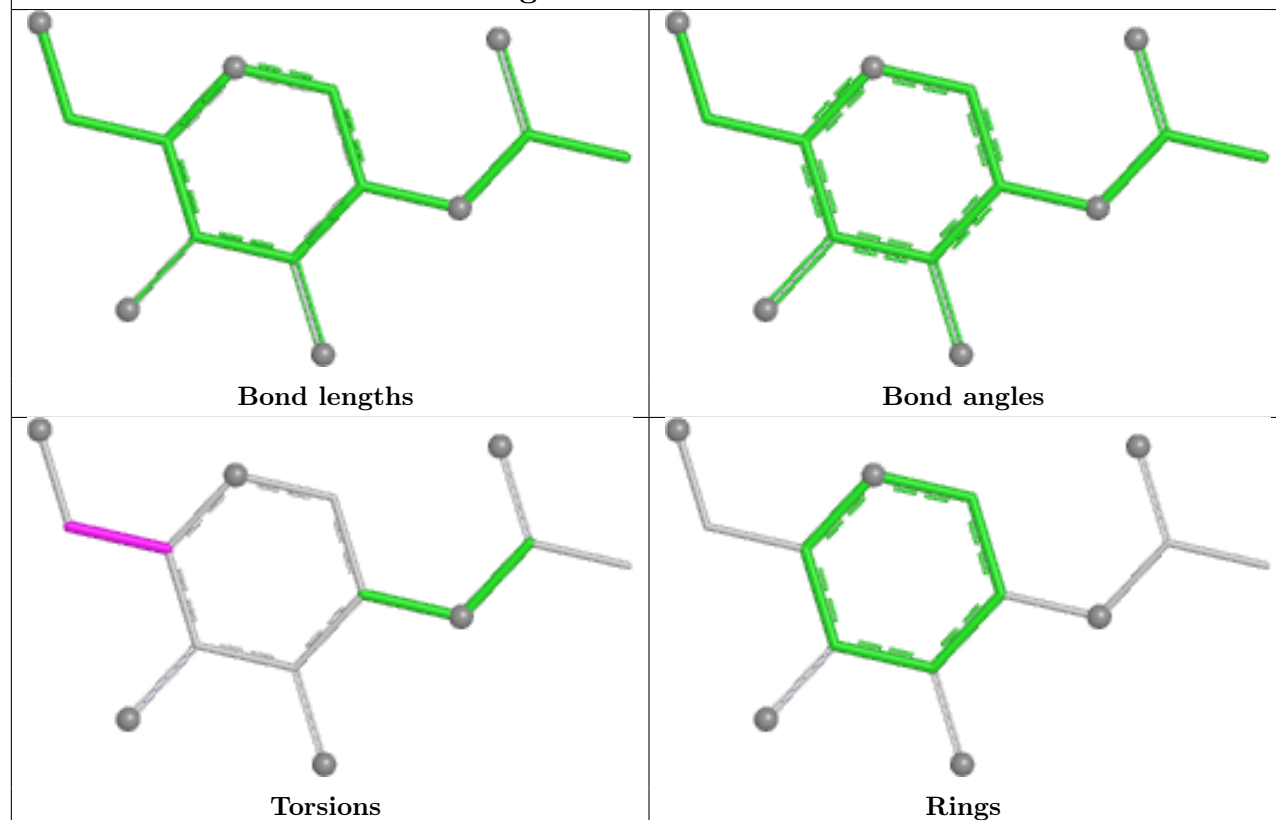
Ligand NAG B 1408



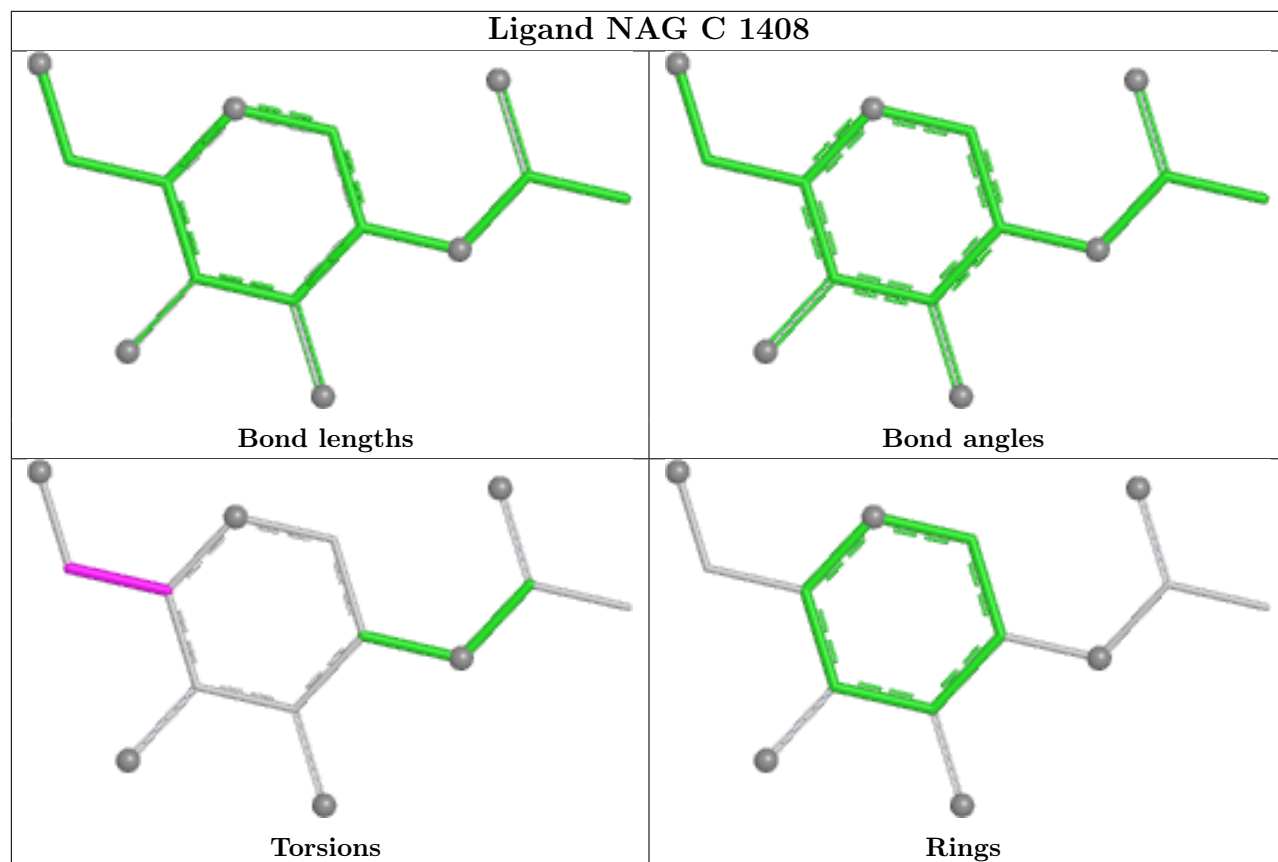
Ligand NAG A 1409



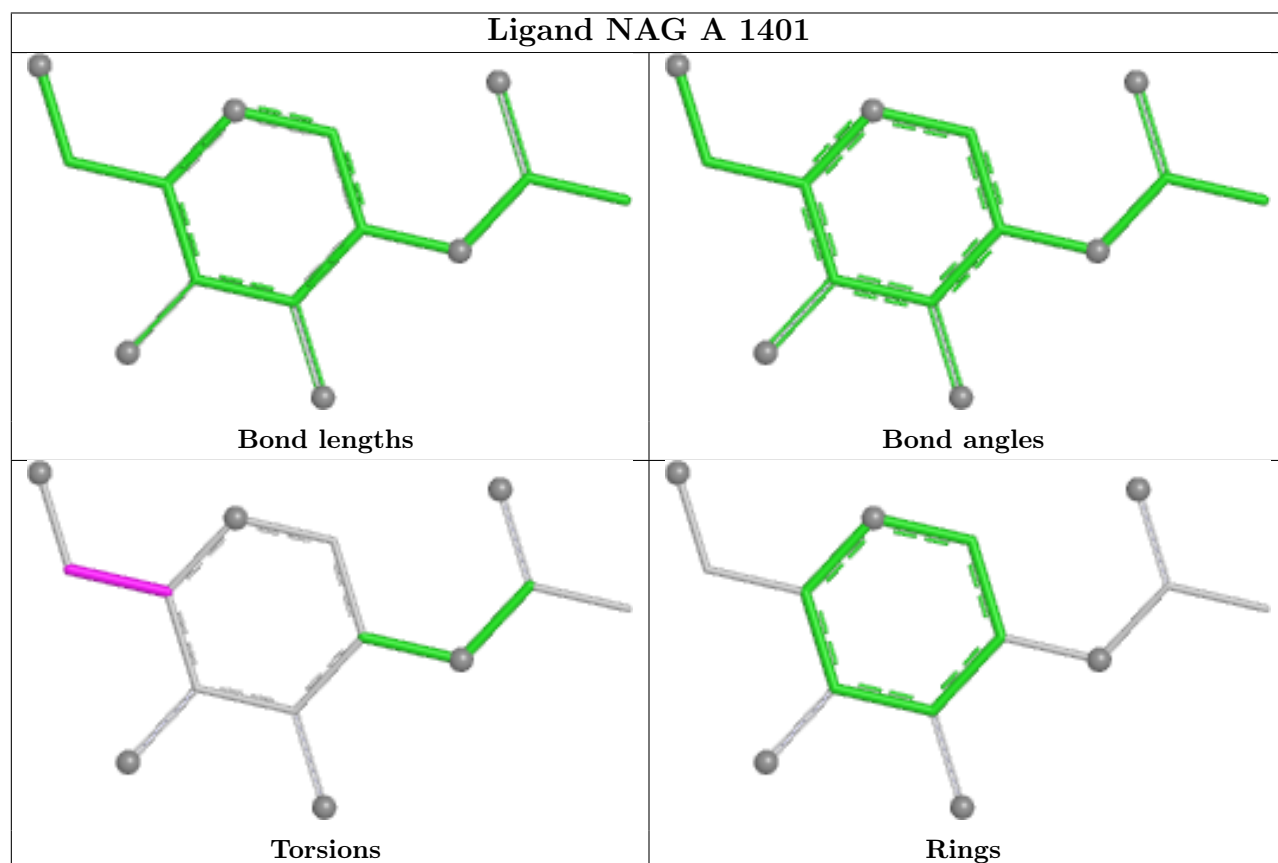
Ligand NAG B 1403



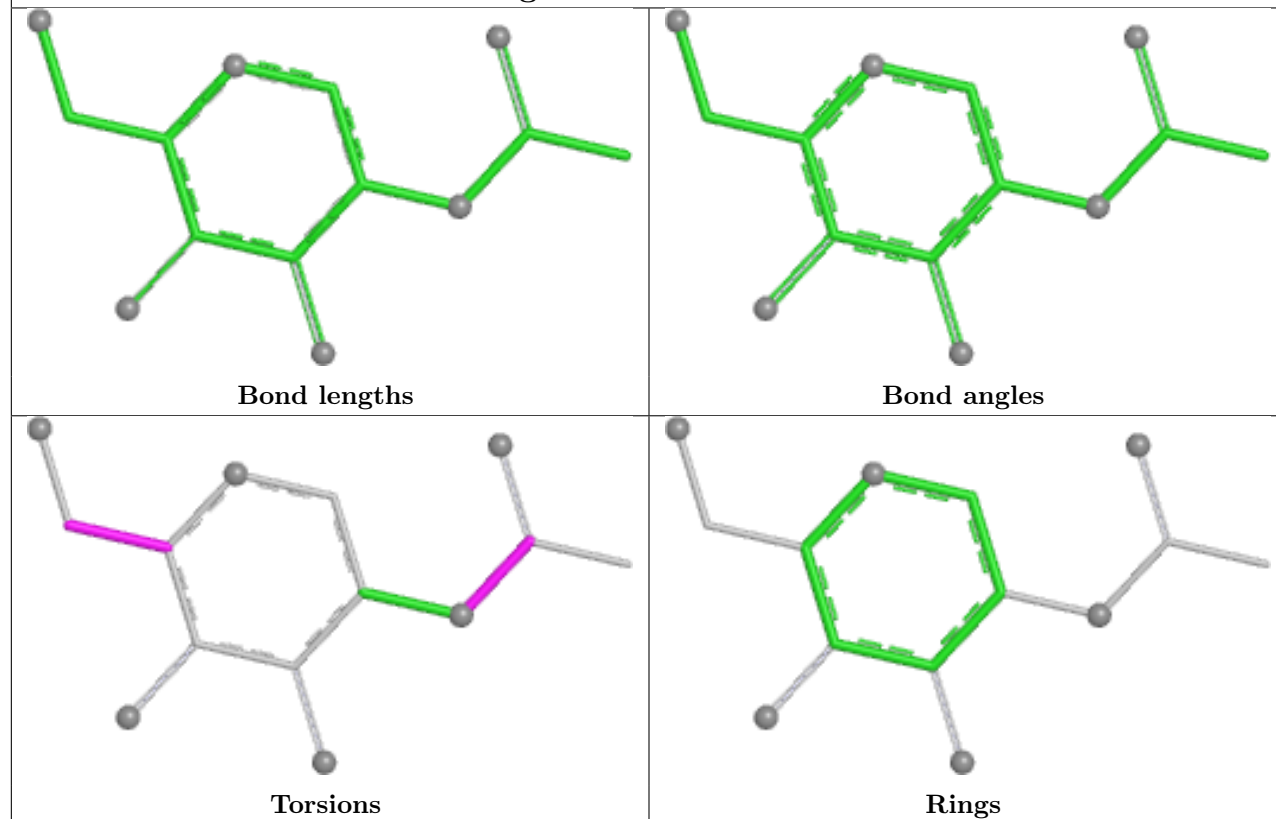
Ligand NAG C 1408



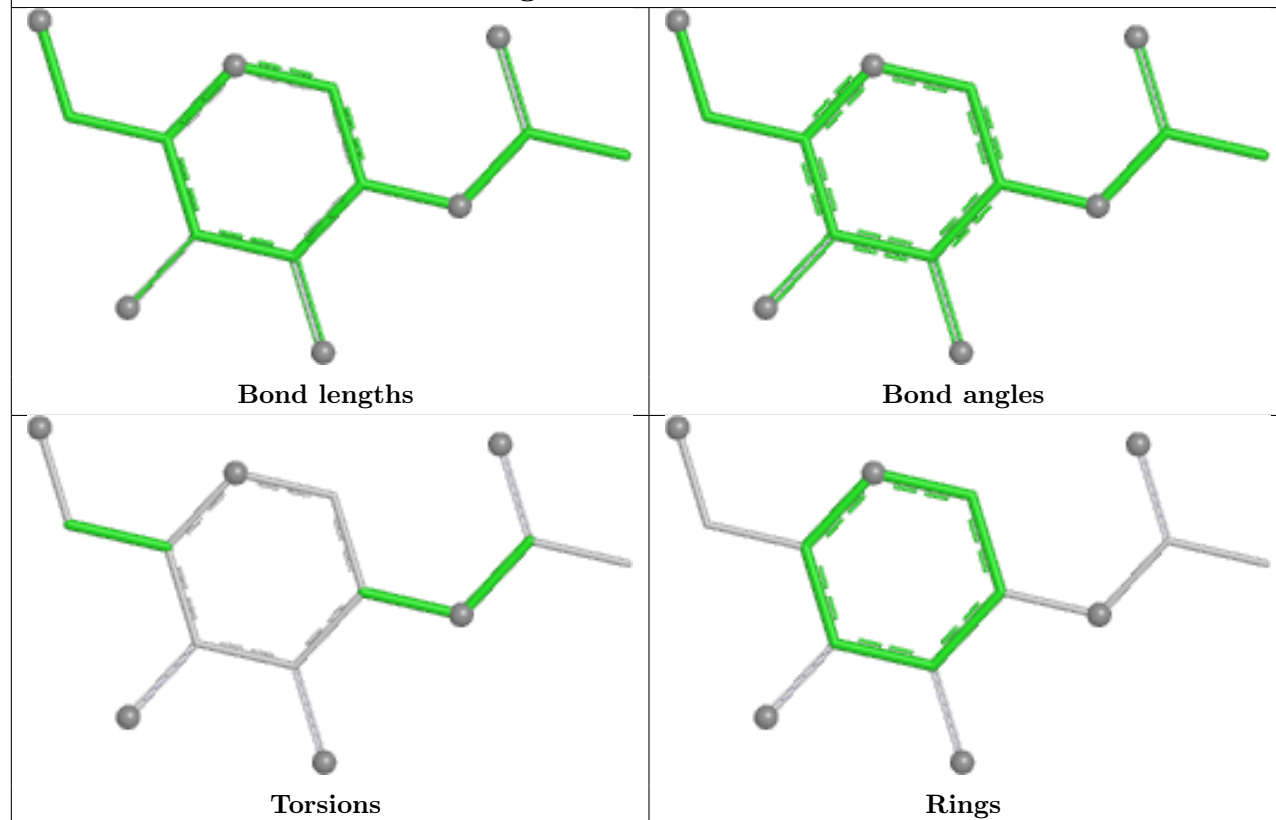
Ligand NAG A 1401



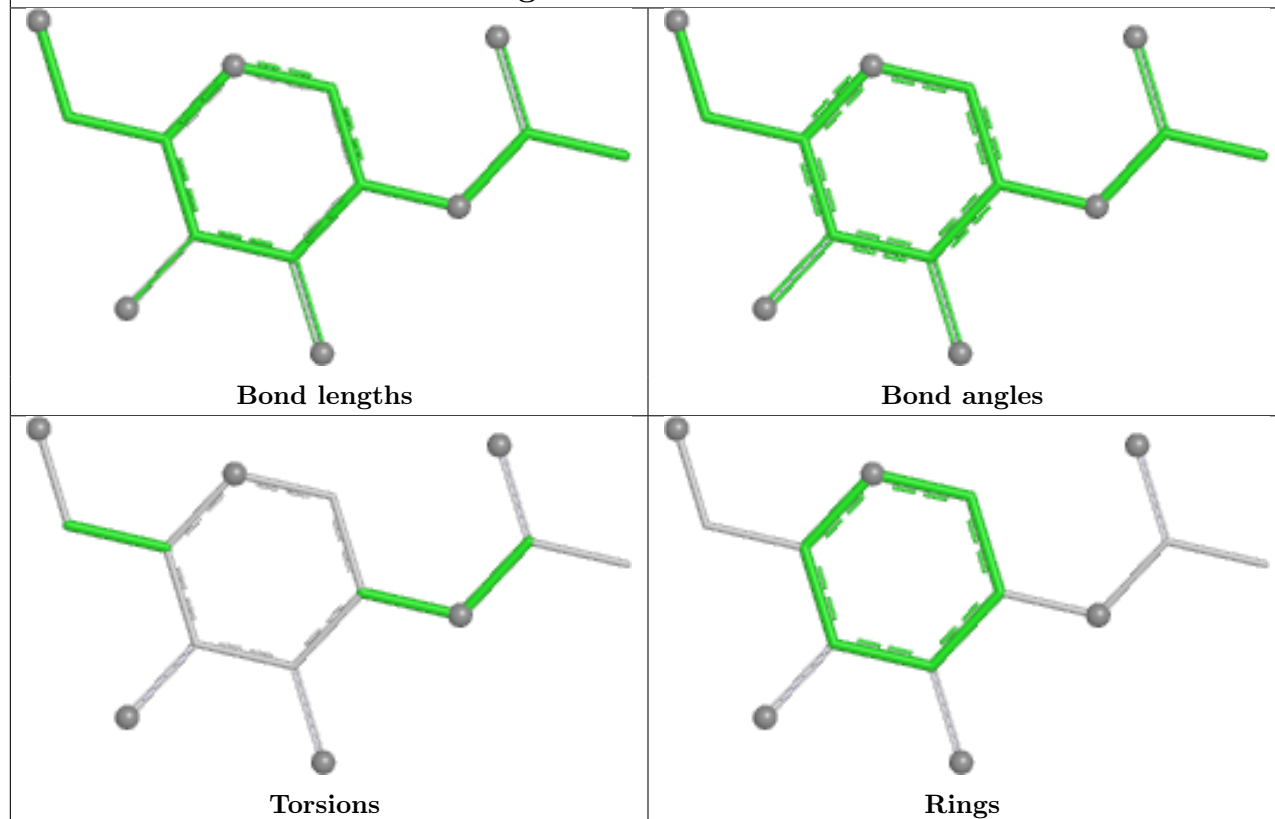
Ligand NAG C 1402



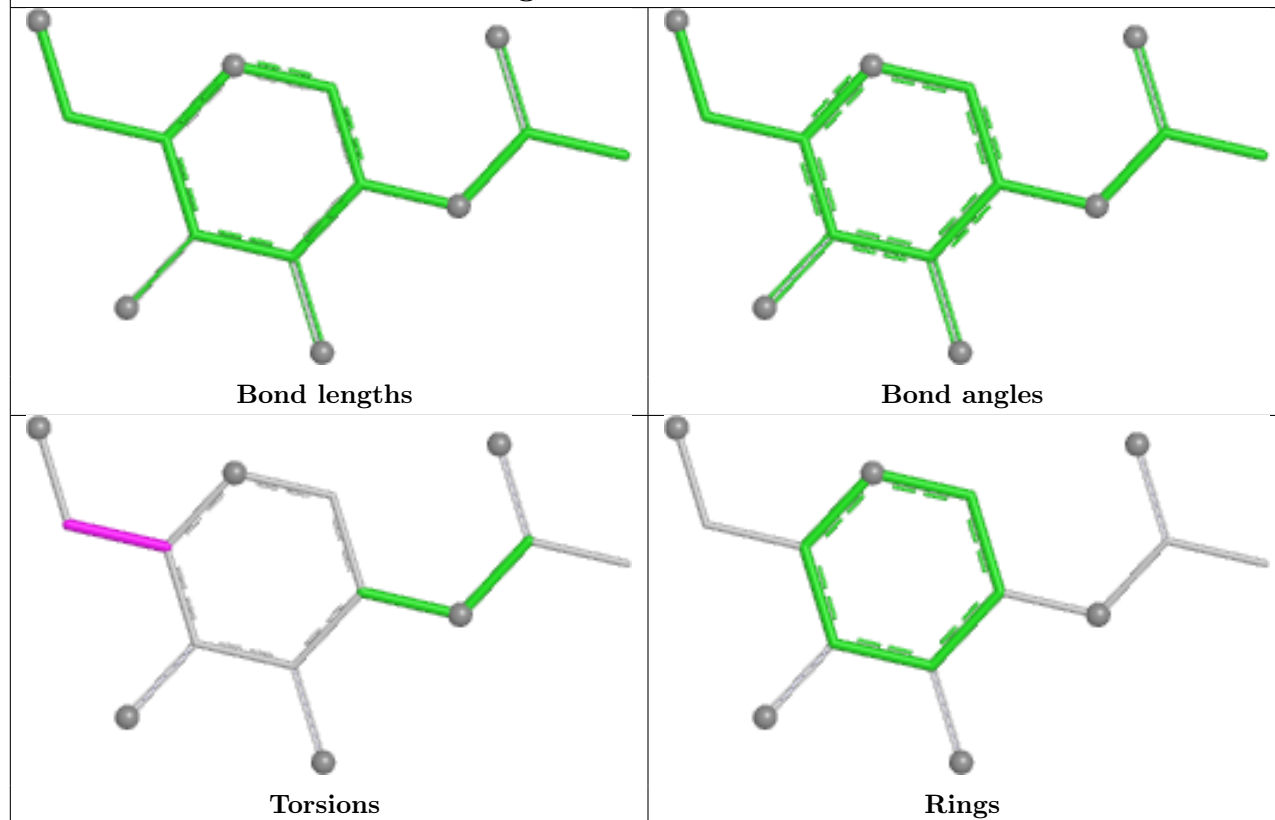
Ligand NAG A 1403

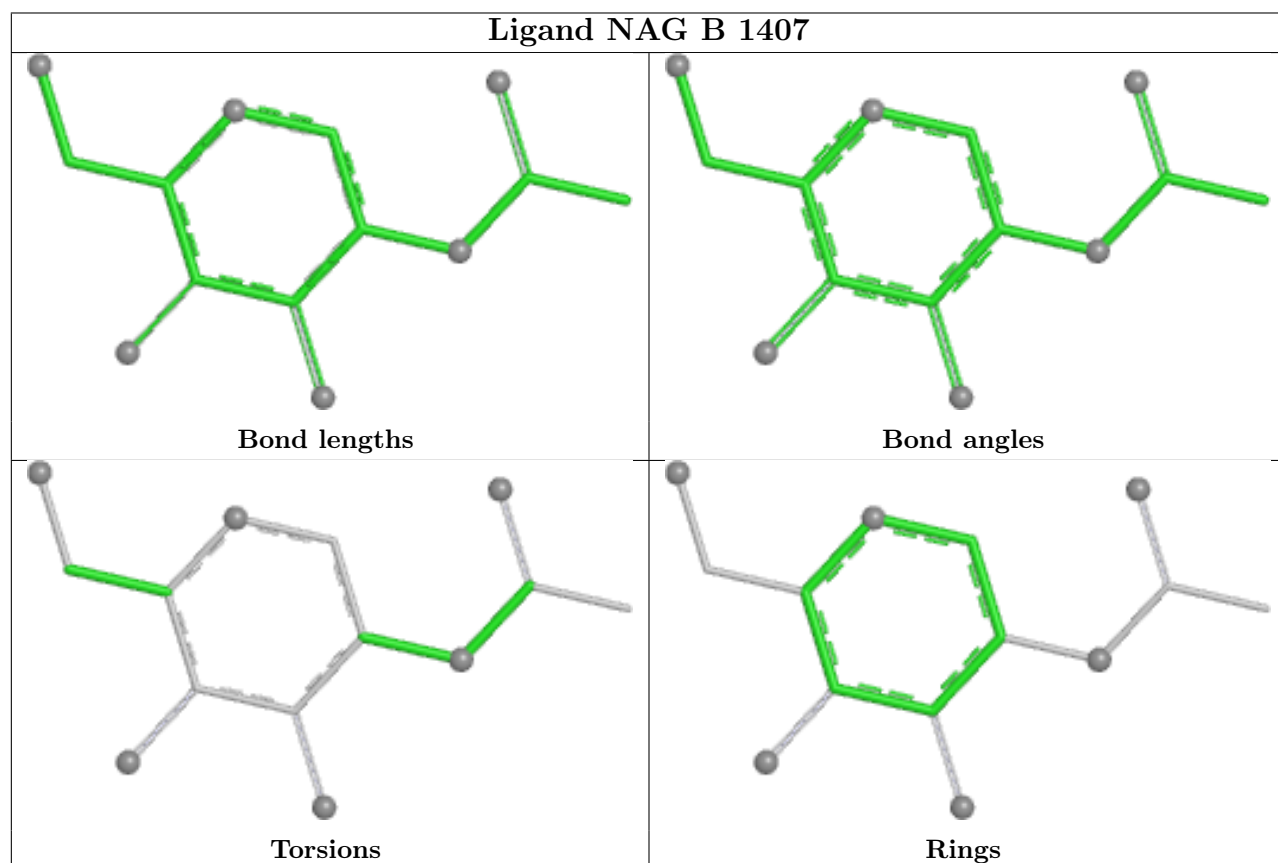
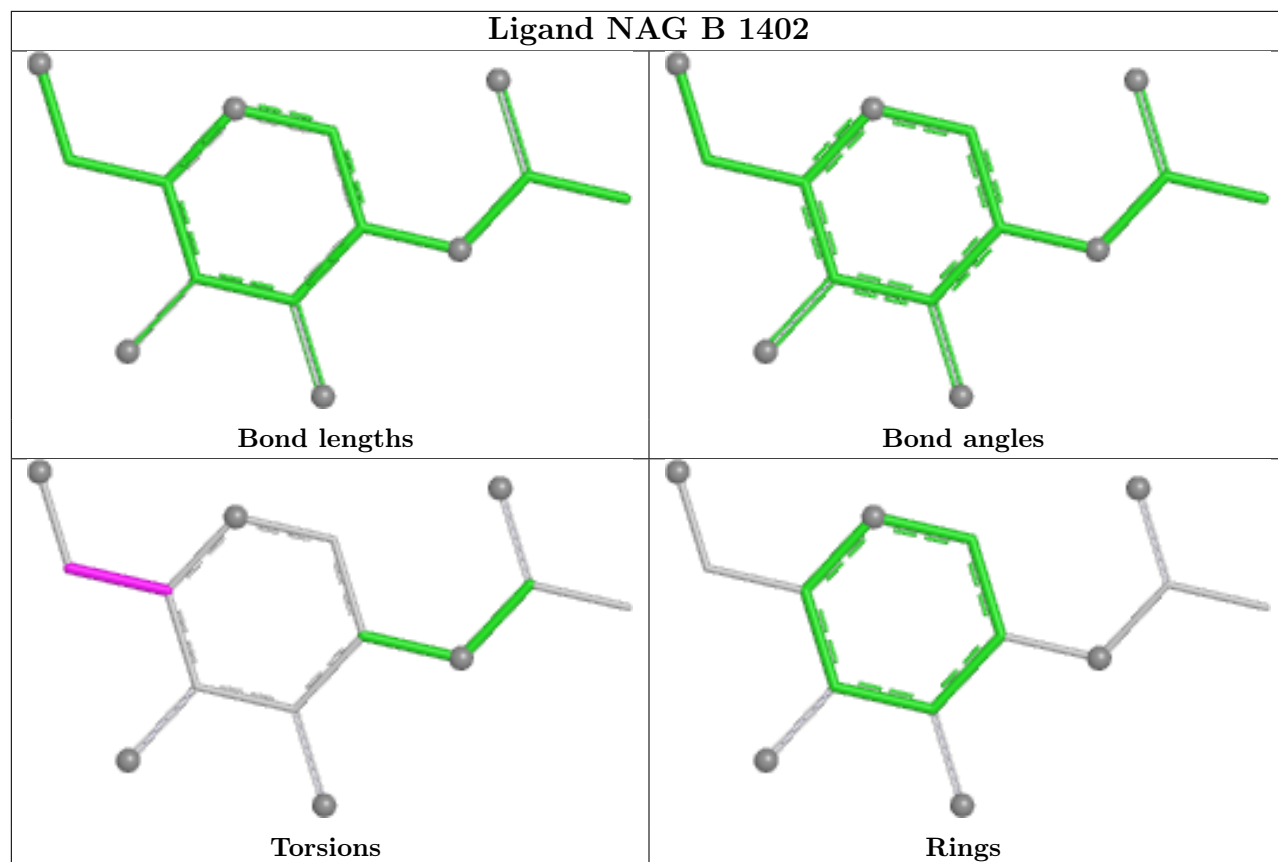


Ligand NAG C 1404

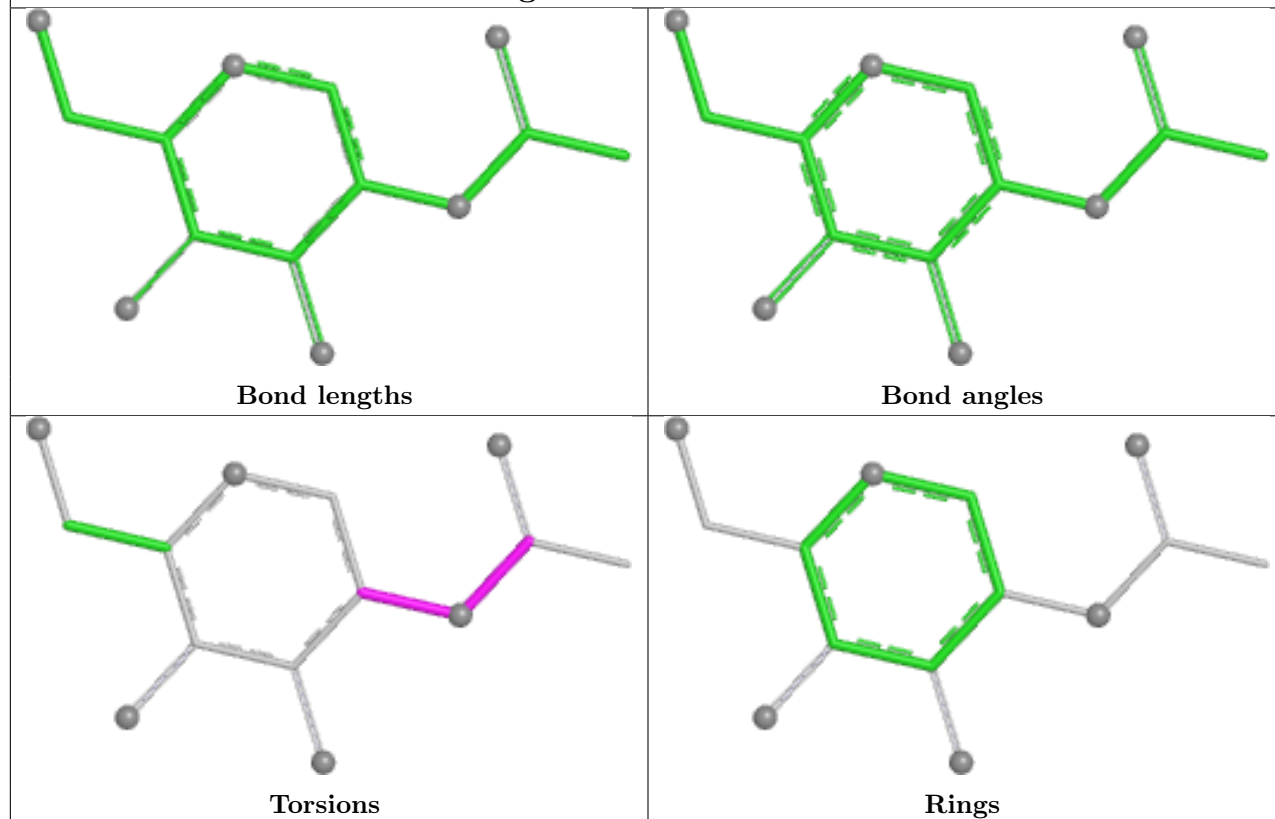


Ligand NAG A 1404

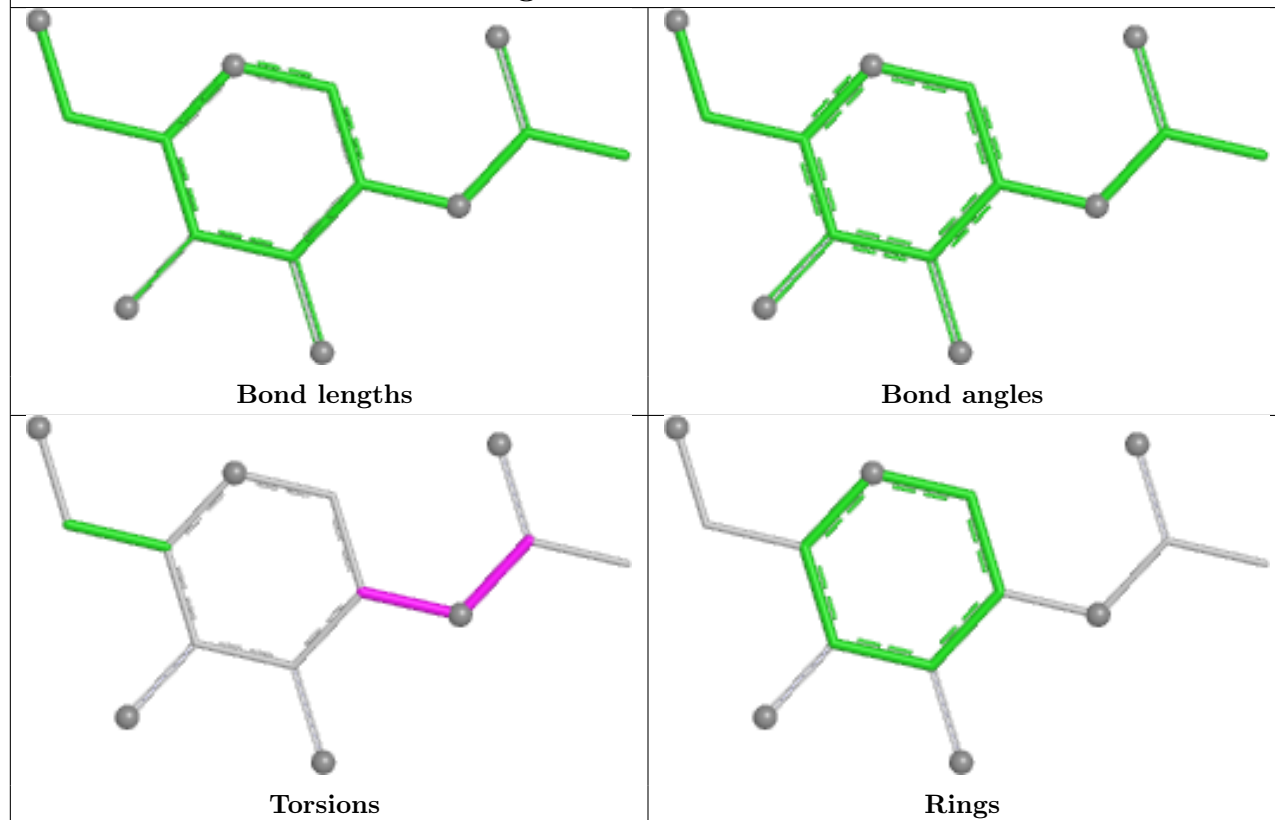




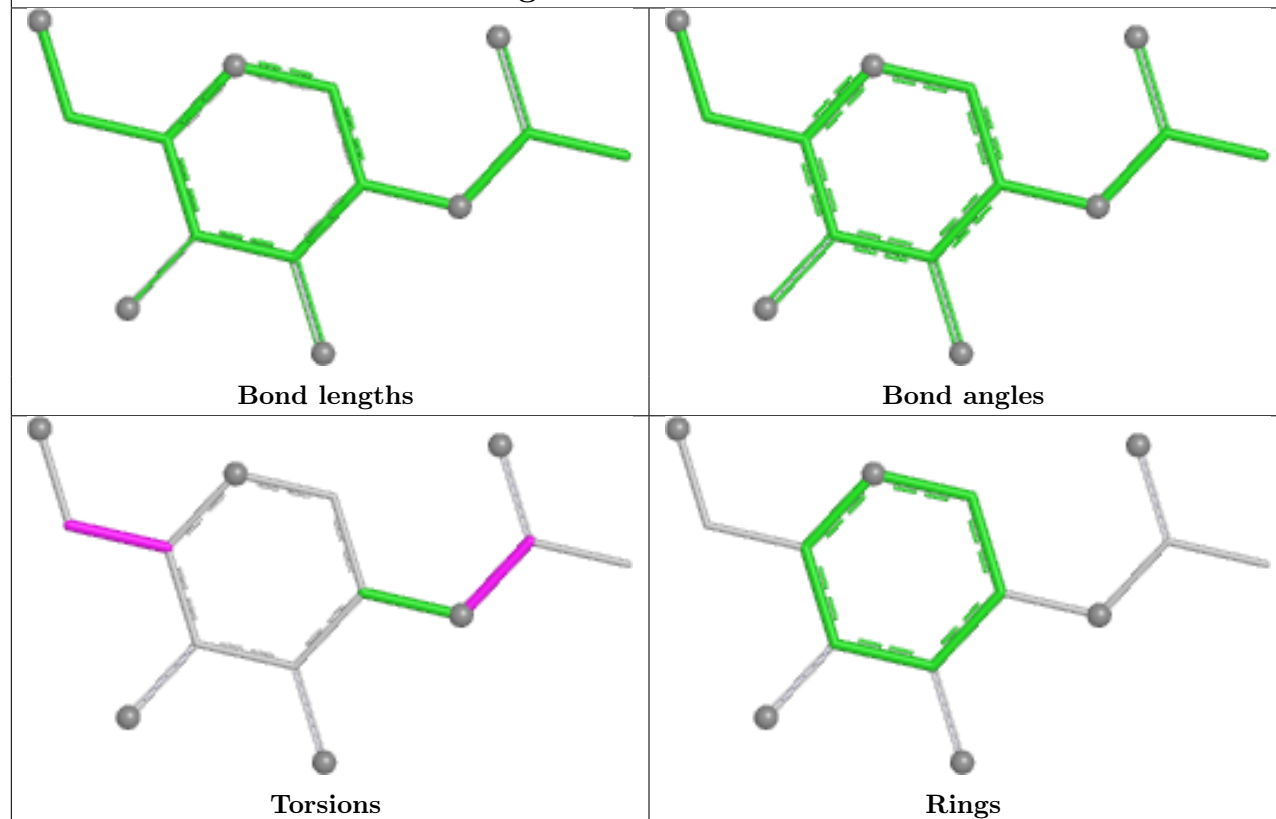
Ligand NAG B 1405



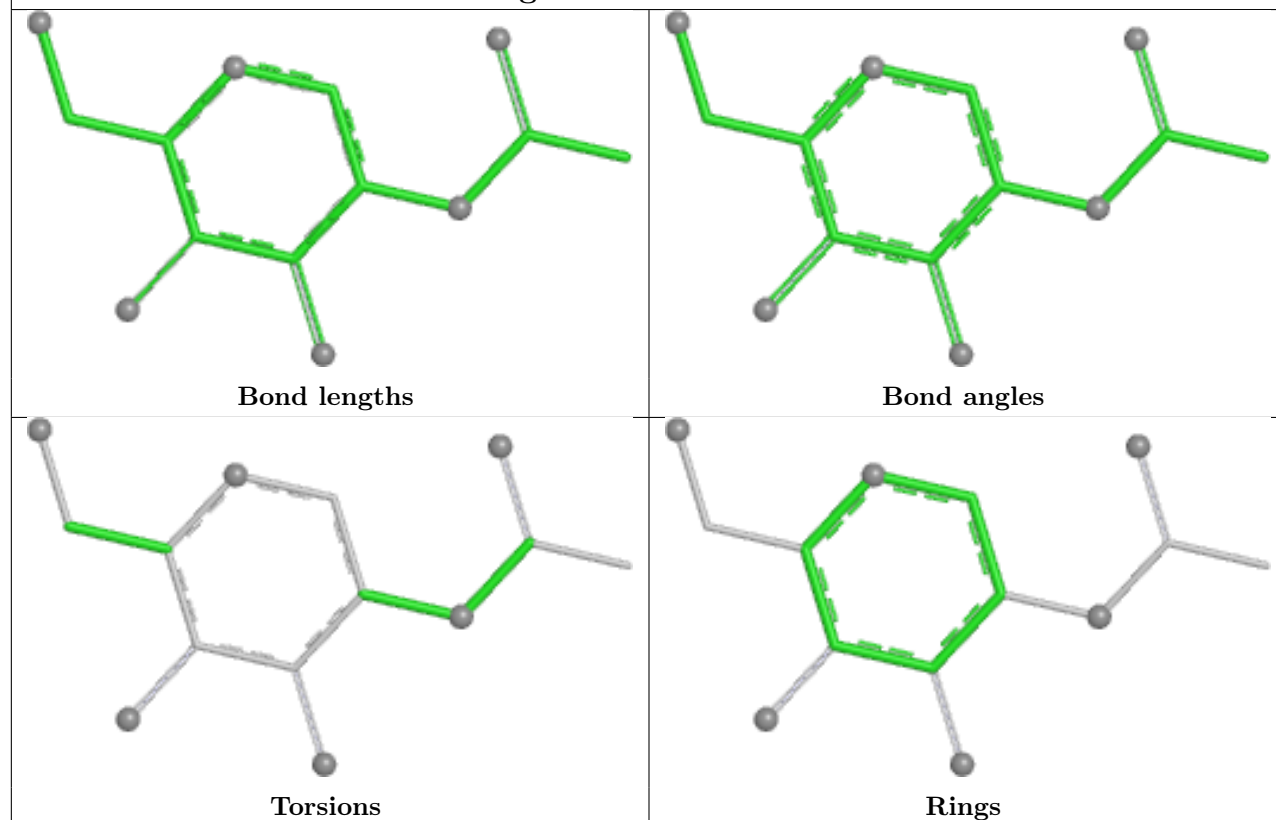
Ligand NAG A 1406



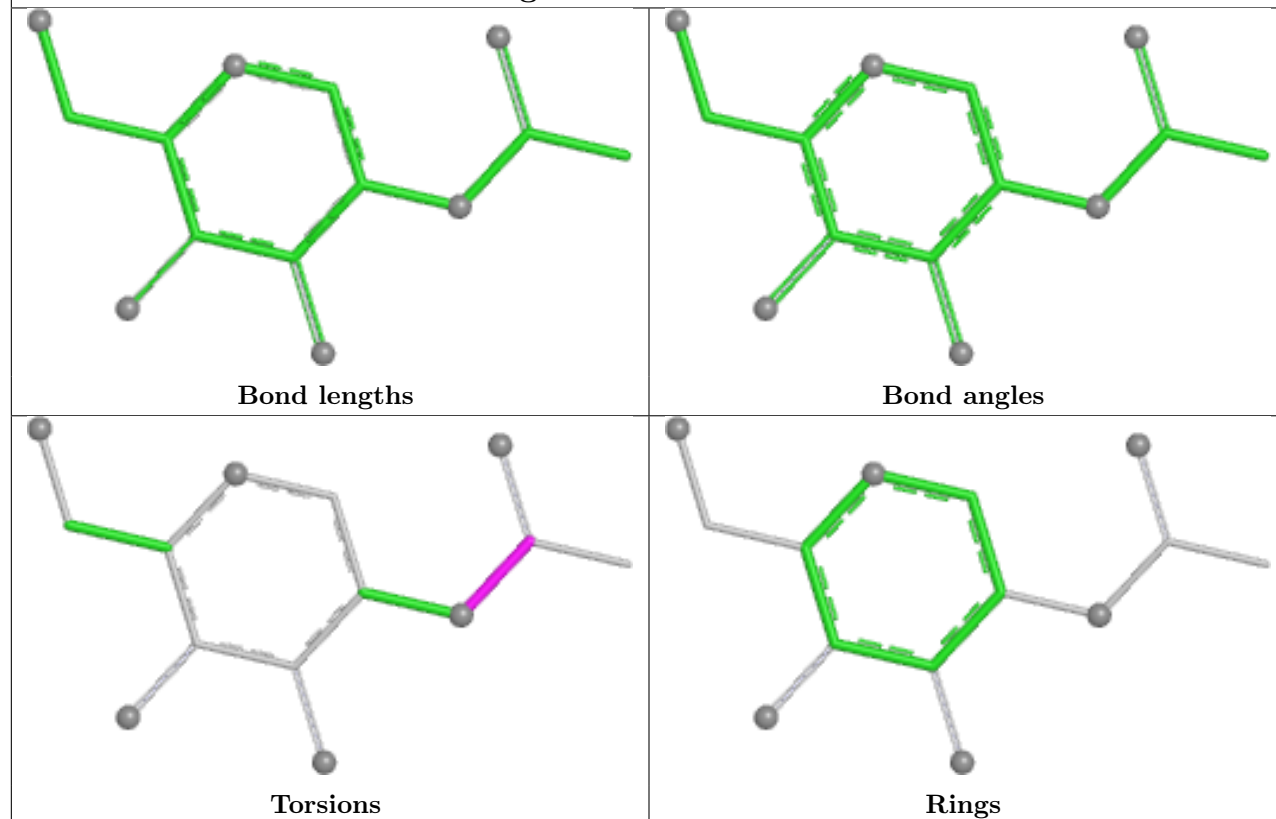
Ligand NAG A 1402



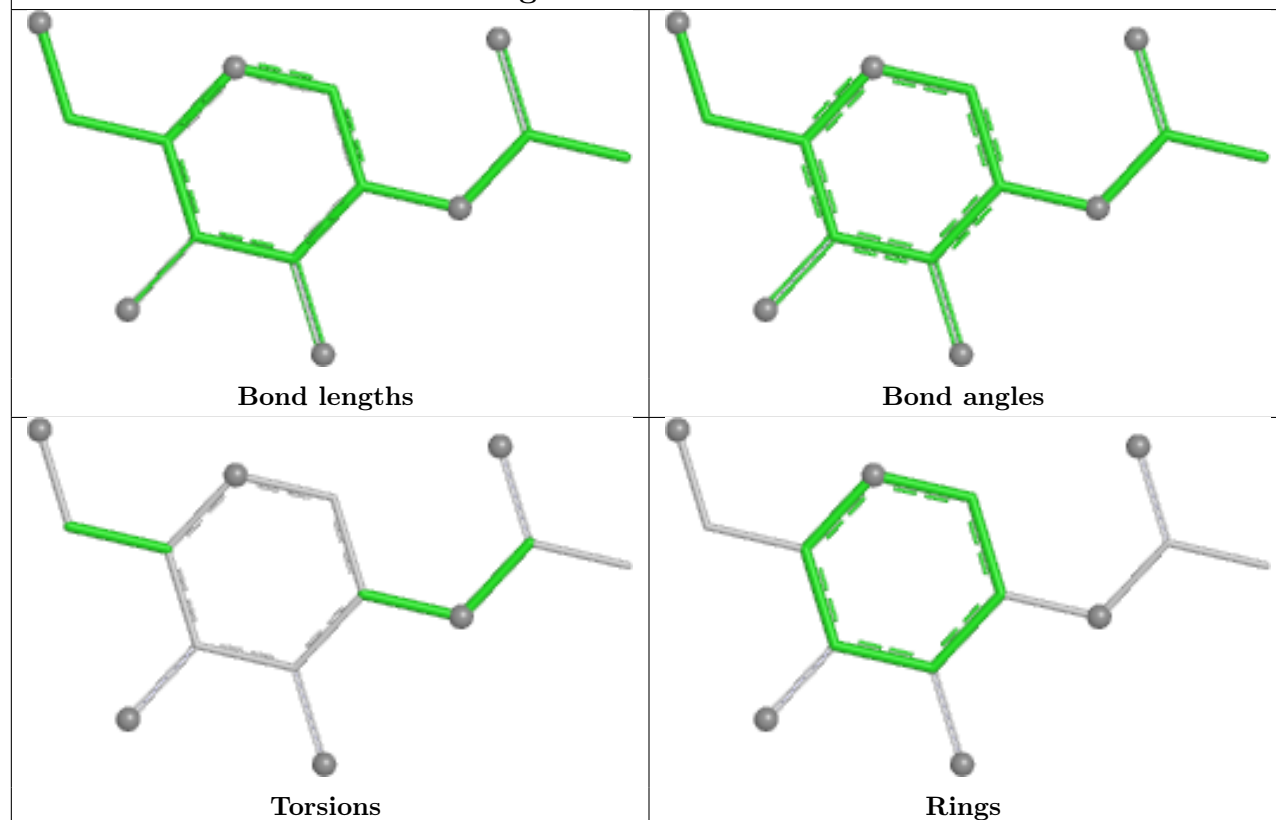
Ligand NAG C 1406



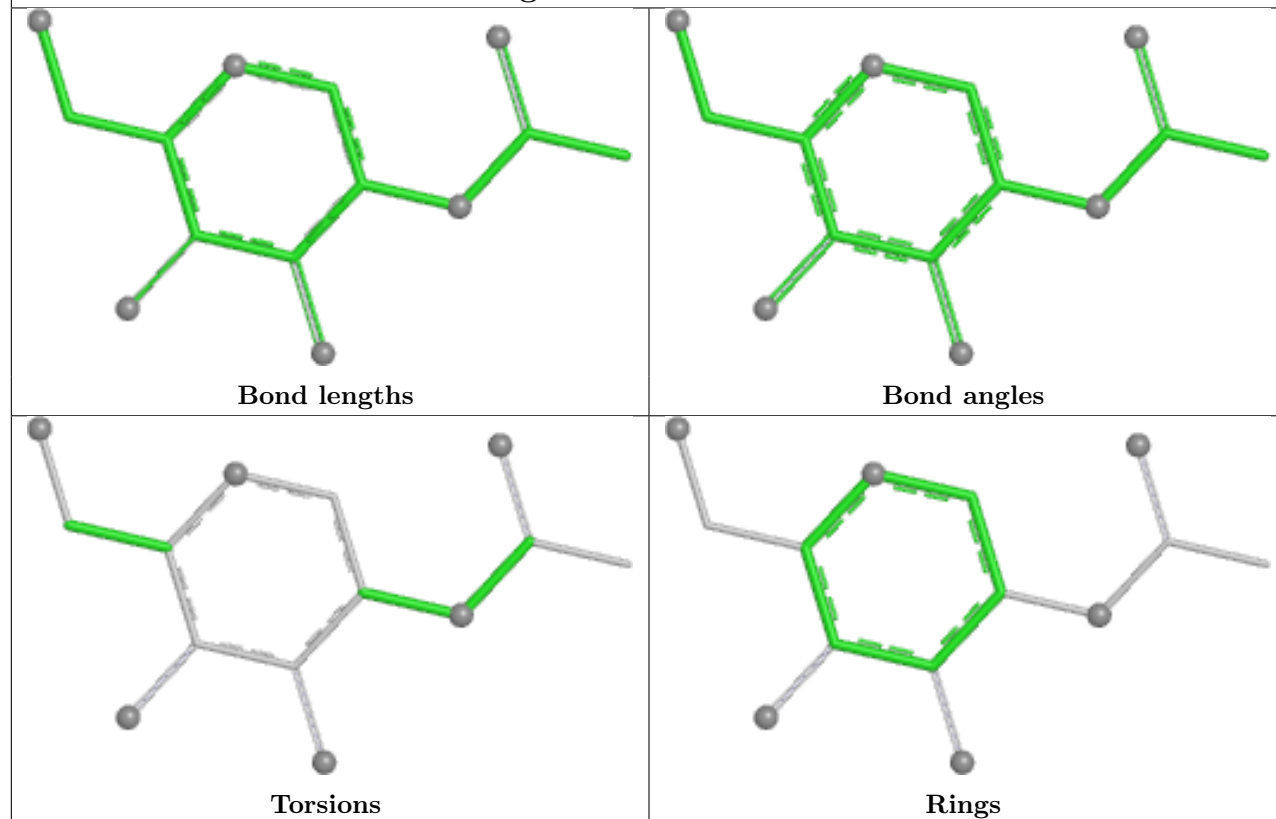
Ligand NAG B 1401



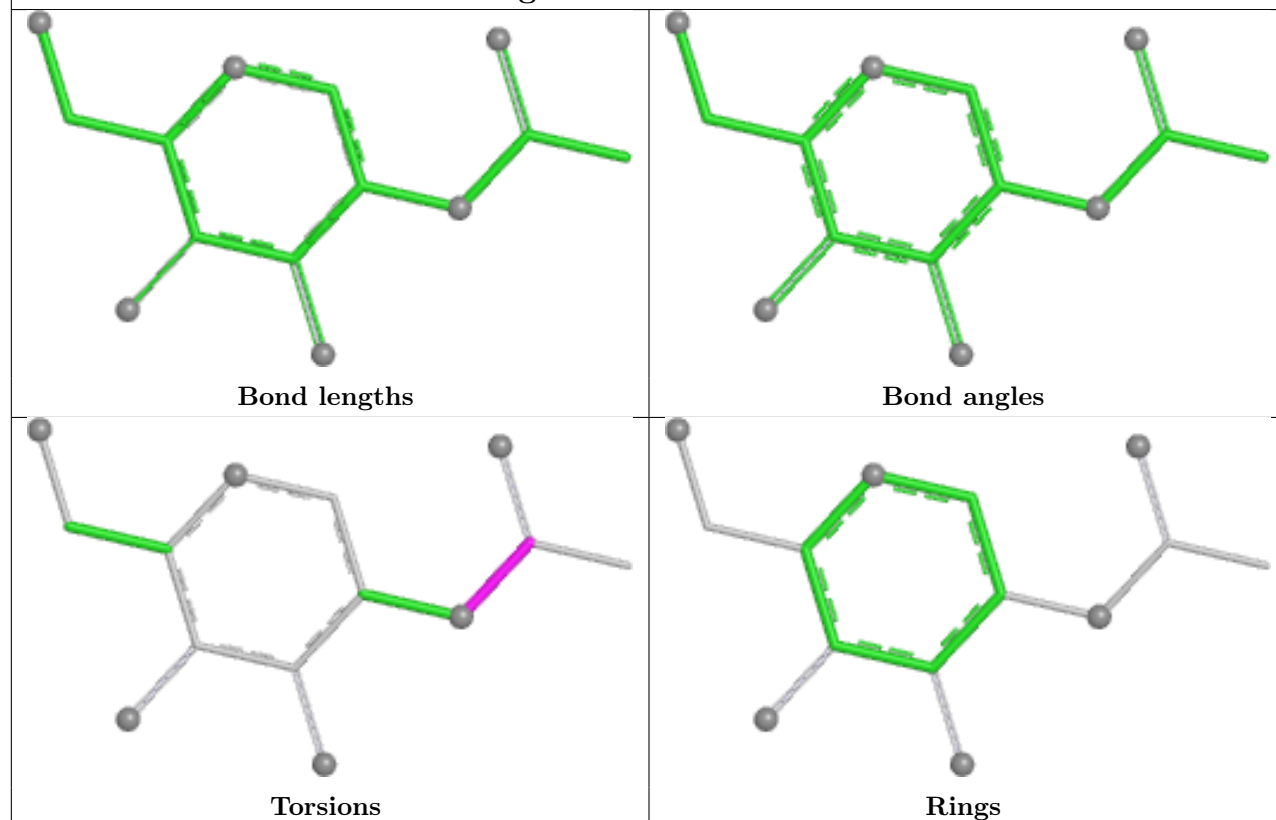
Ligand NAG C 1401



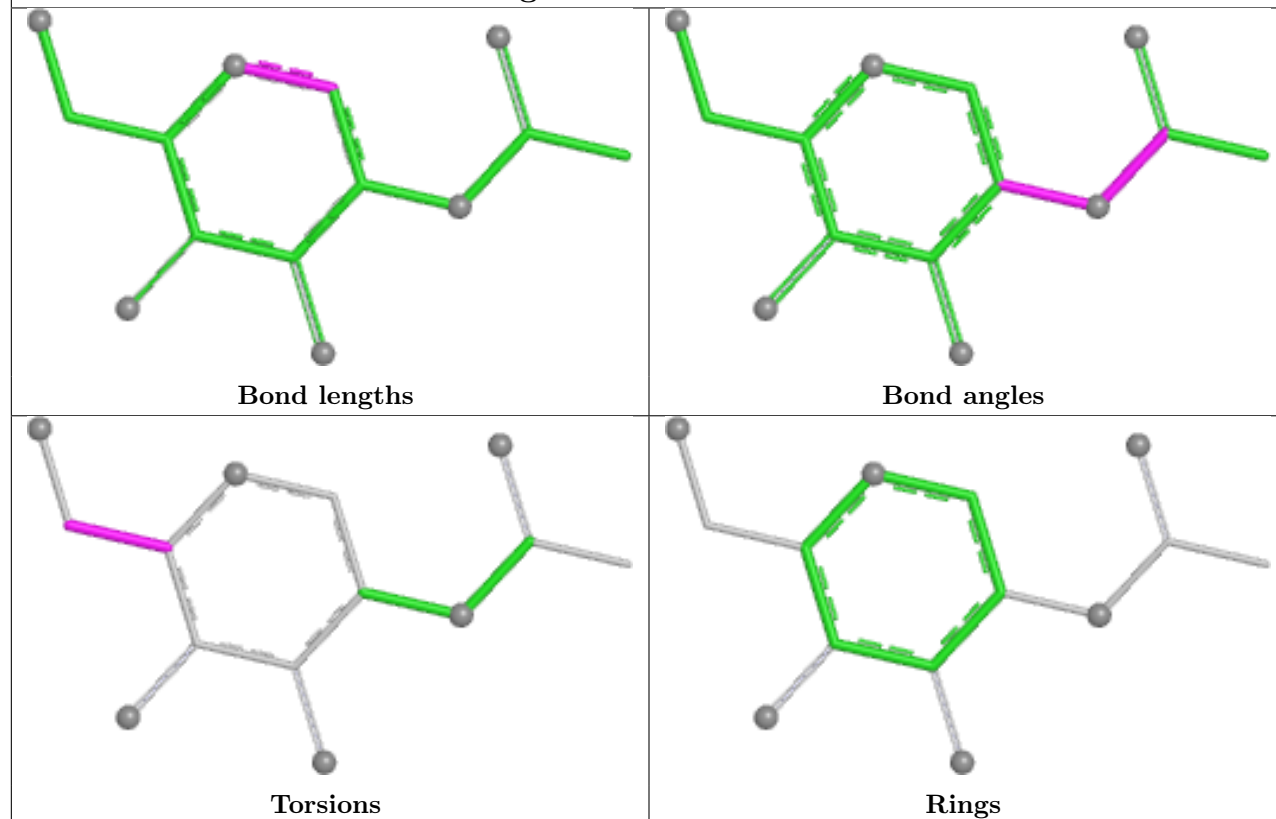
Ligand NAG B 1406



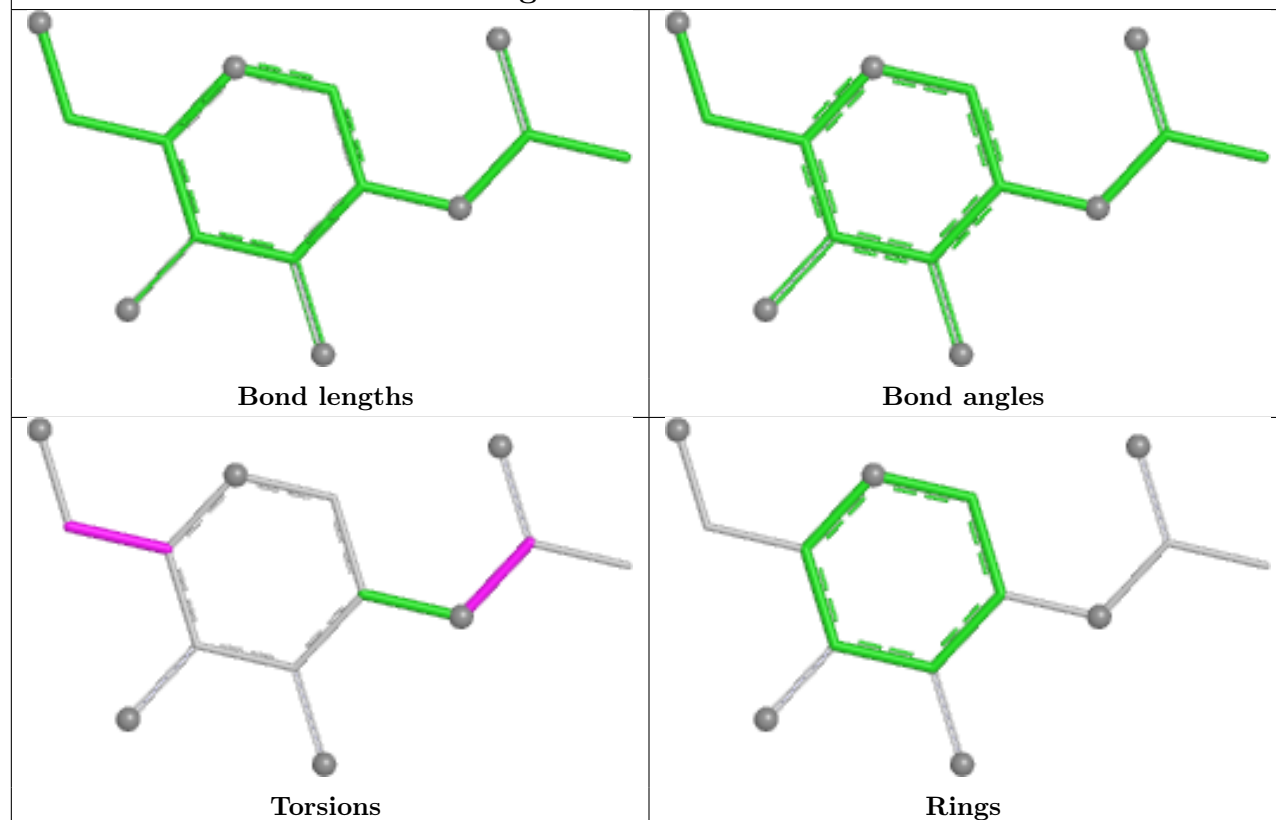
Ligand NAG A 1405



Ligand NAG B 1404



Ligand NAG C 1405



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.

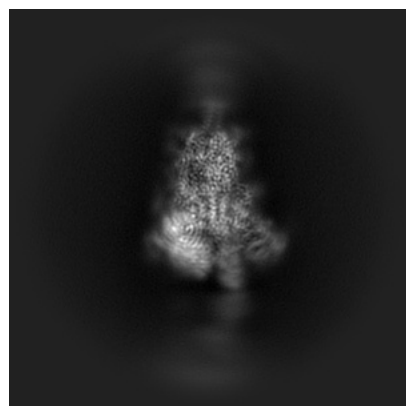
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-40094. These allow visual inspection of the internal detail of the map and identification of artifacts.

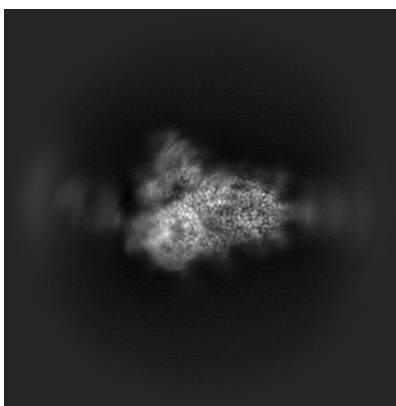
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

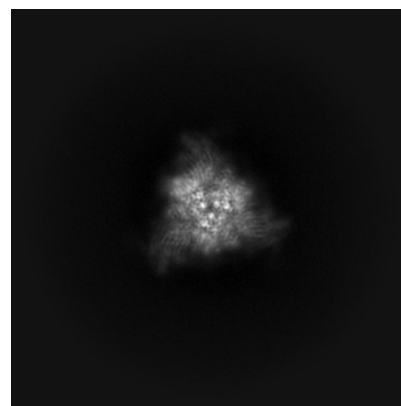
6.1.1 Primary map



X

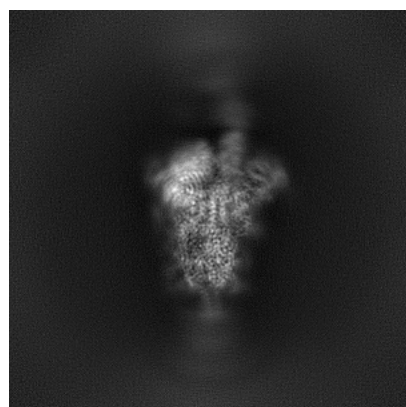


Y

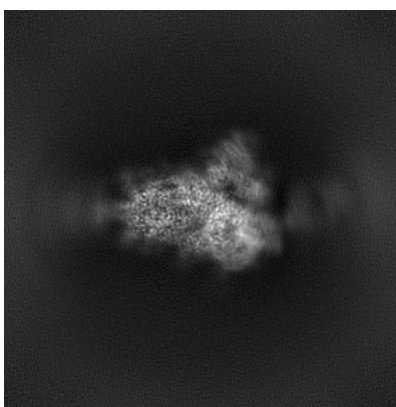


Z

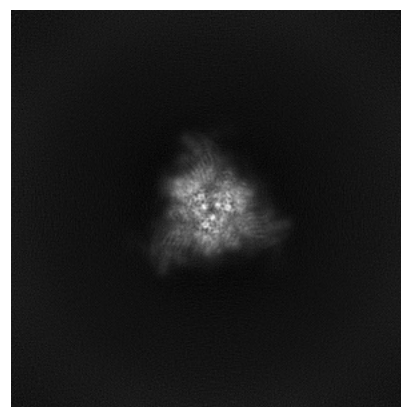
6.1.2 Raw map



X



Y

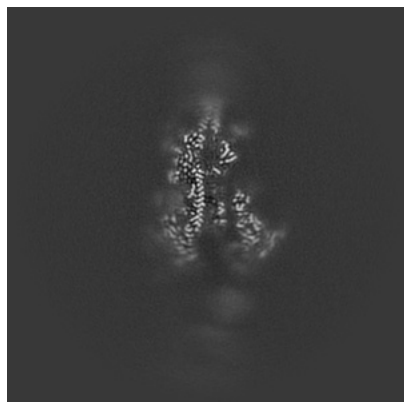


Z

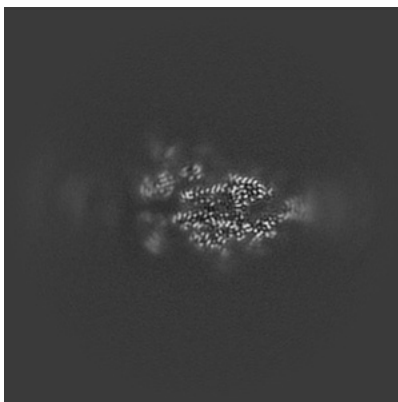
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

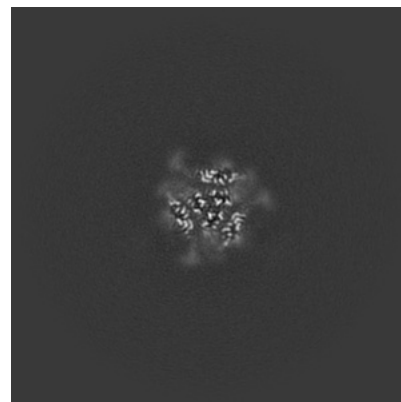
6.2.1 Primary map



X Index: 240

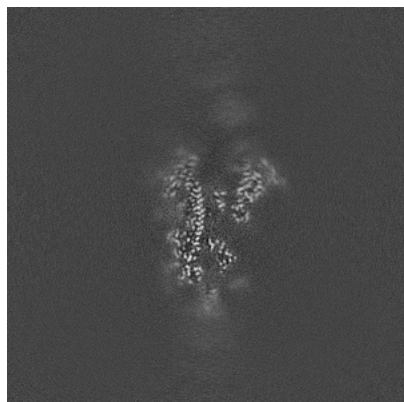


Y Index: 240

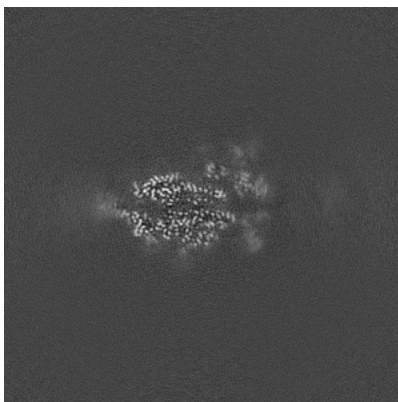


Z Index: 240

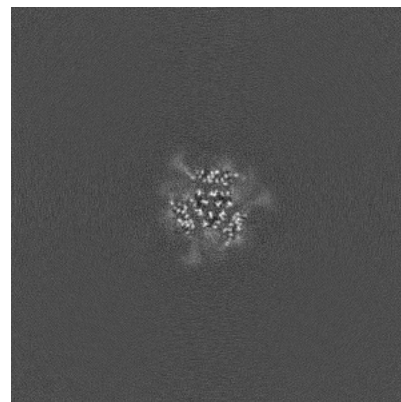
6.2.2 Raw map



X Index: 240



Y Index: 240

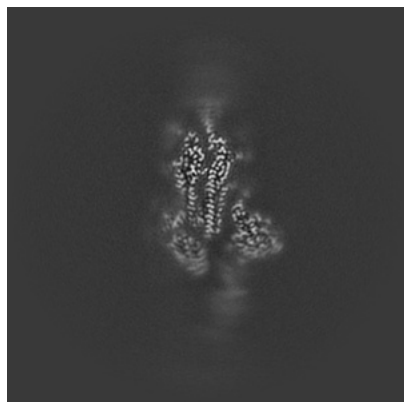


Z Index: 240

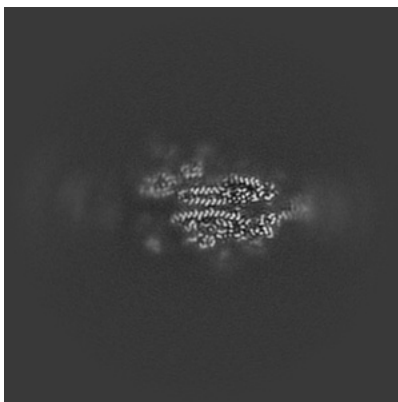
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

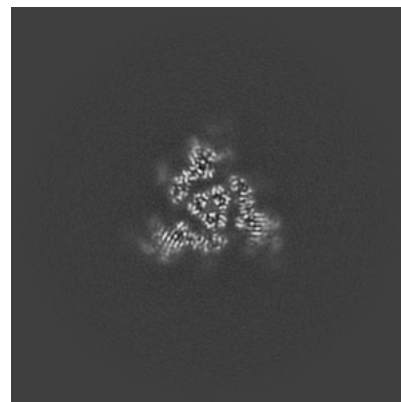
6.3.1 Primary map



X Index: 229

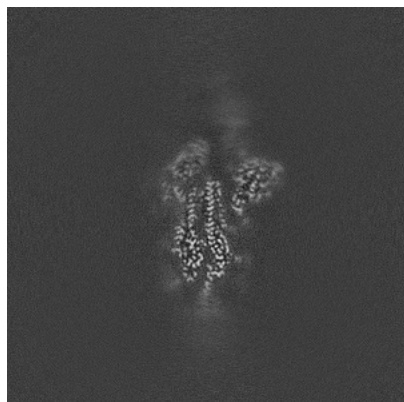


Y Index: 243

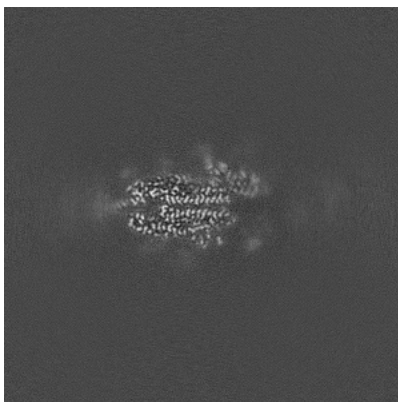


Z Index: 218

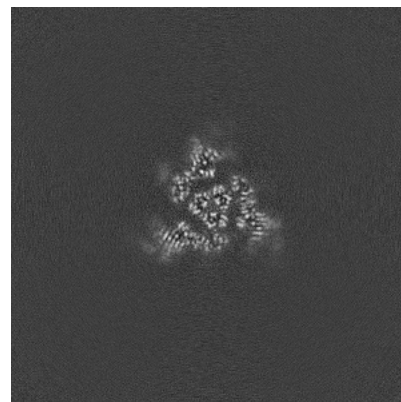
6.3.2 Raw map



X Index: 231



Y Index: 246

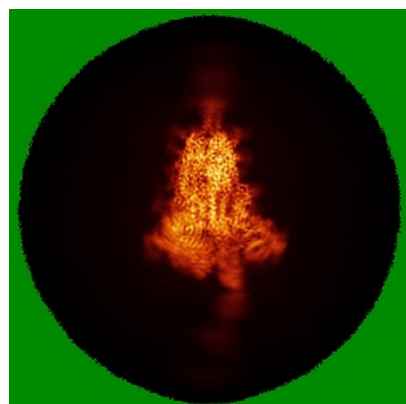


Z Index: 261

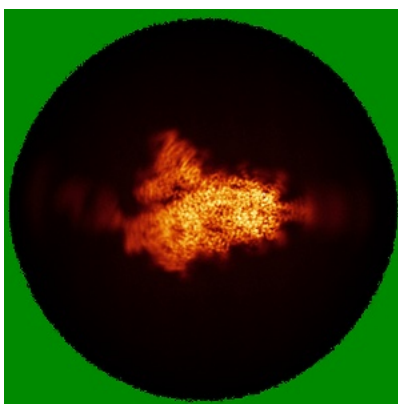
The images above show the largest variance slices of the map in three orthogonal directions.

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

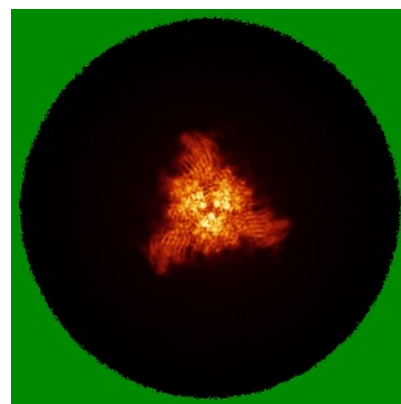
6.4.1 Primary map



X



Y

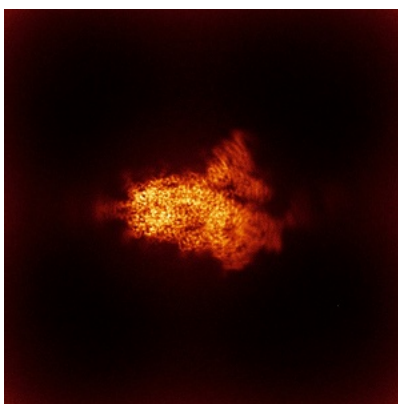


Z

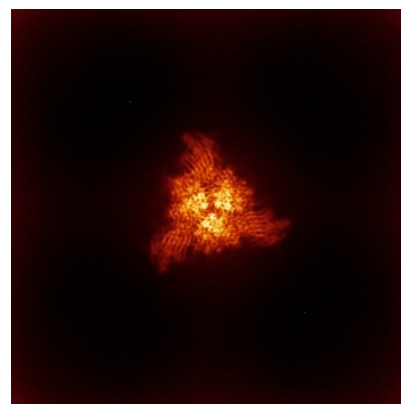
6.4.2 Raw map



X



Y

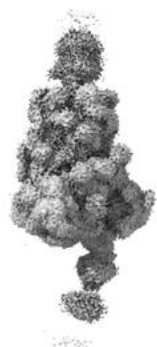


Z

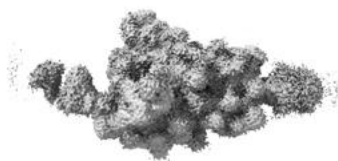
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



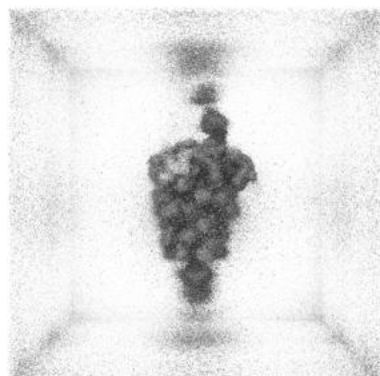
Y



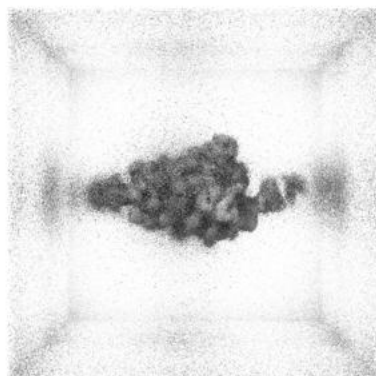
Z

The images above show the 3D surface view of the map at the recommended contour level 0.02. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

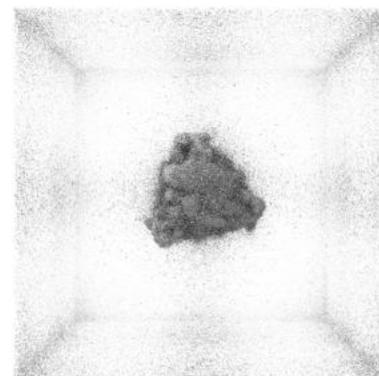
6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

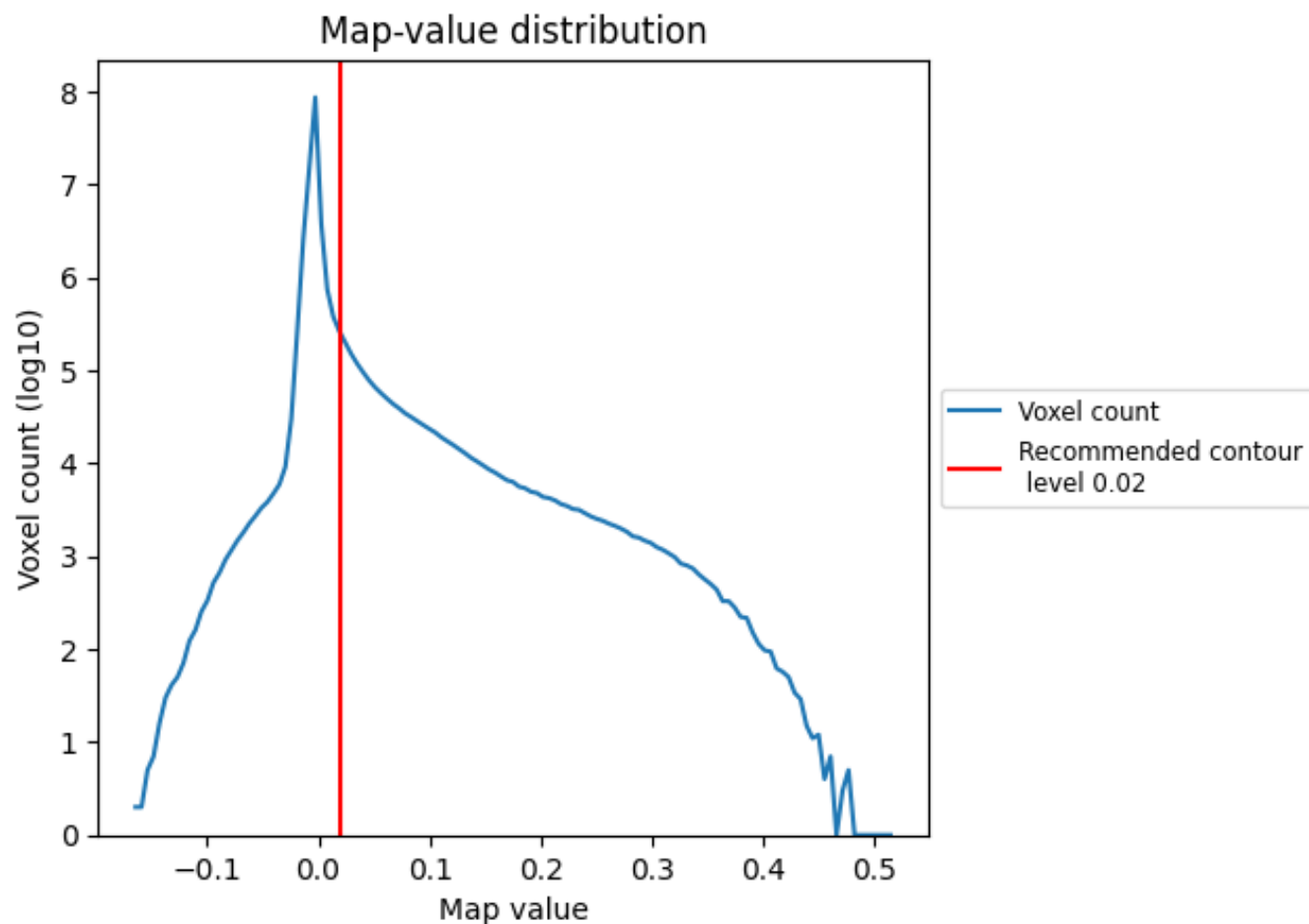
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

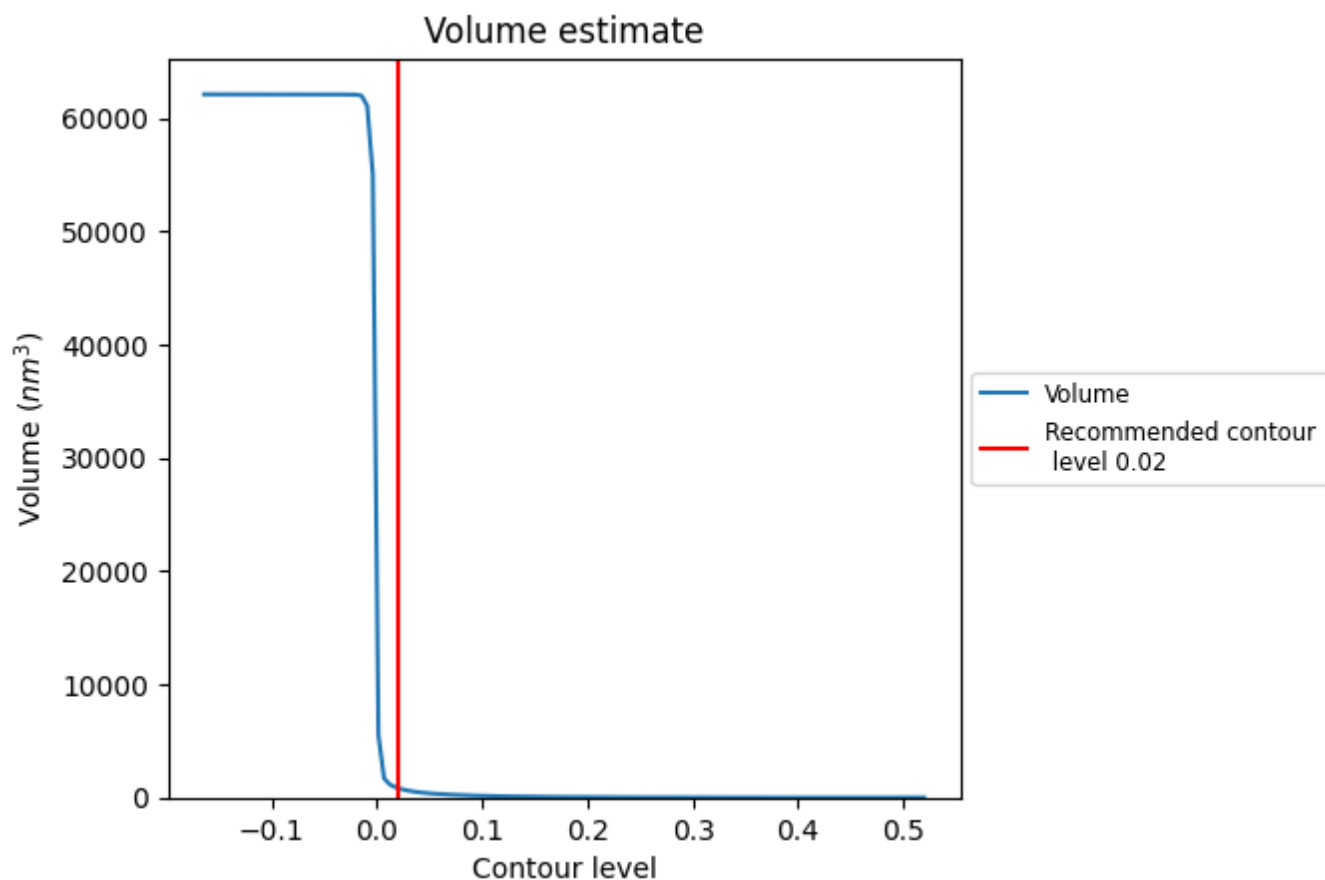
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

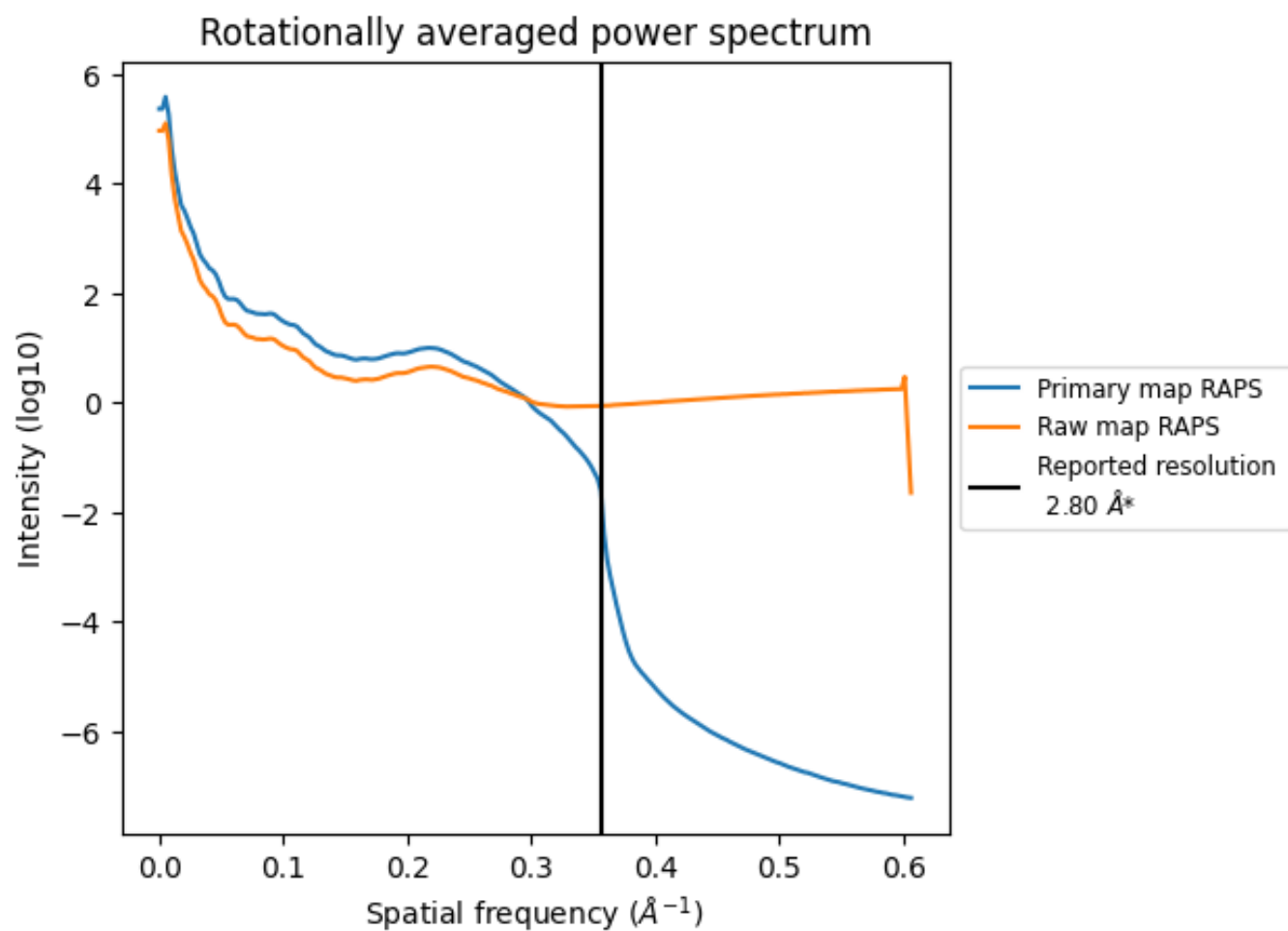
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 838 nm^3 ; this corresponds to an approximate mass of 757 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 ⓘ

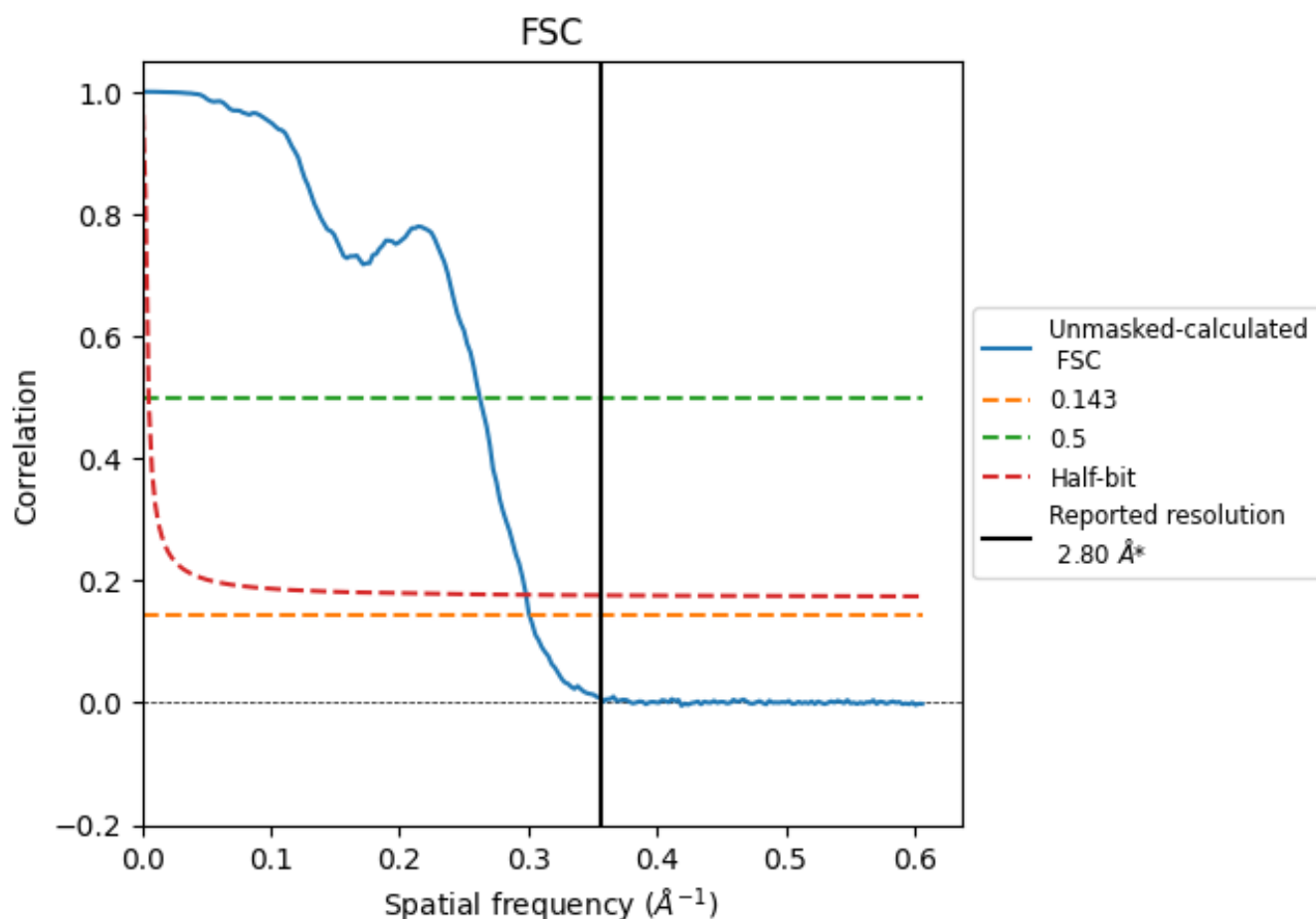


*Reported resolution corresponds to spatial frequency of 0.357 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.357 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.80	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.32	3.81	3.35

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.32 differs from the reported value 2.8 by more than 10 %

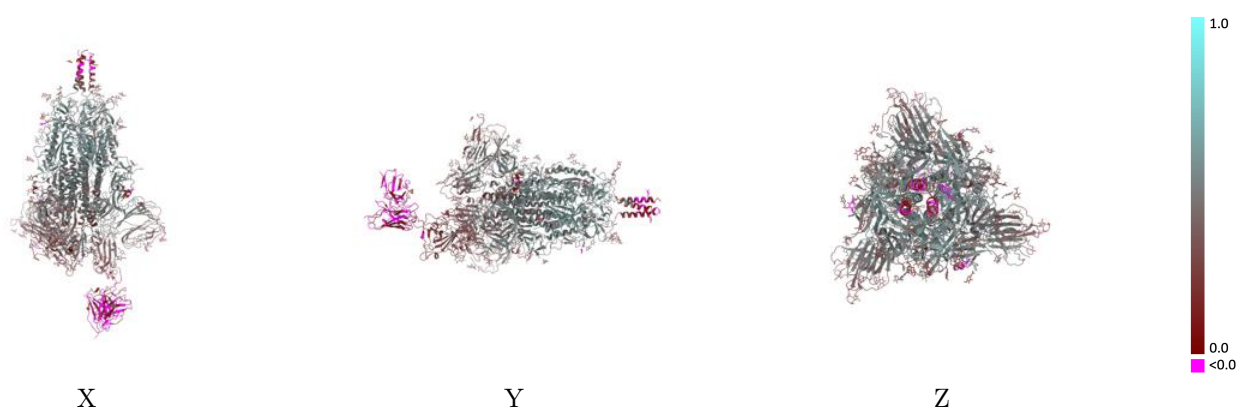
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-40094 and PDB model 8GJM. Per-residue inclusion information can be found in section 3 on page 13.

9.1 Map-model overlay [i](#)

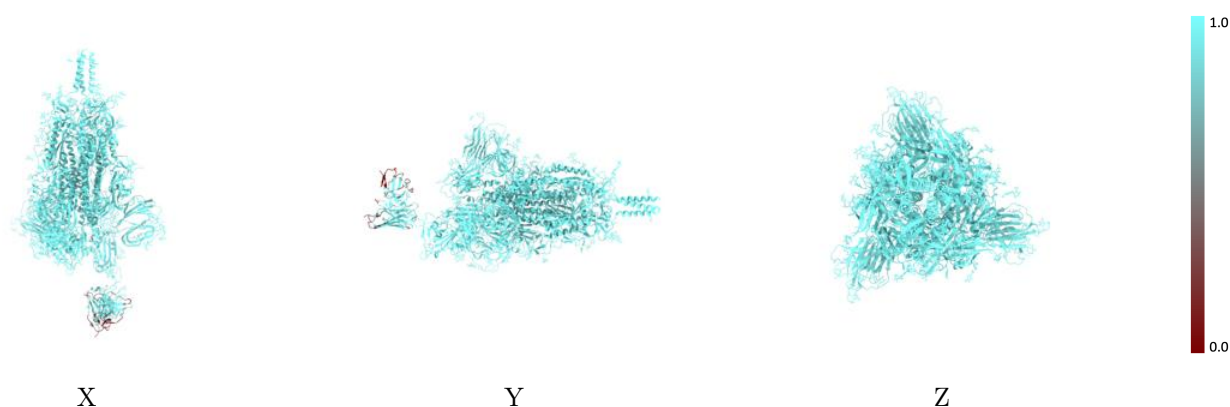
This section was not generated.

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



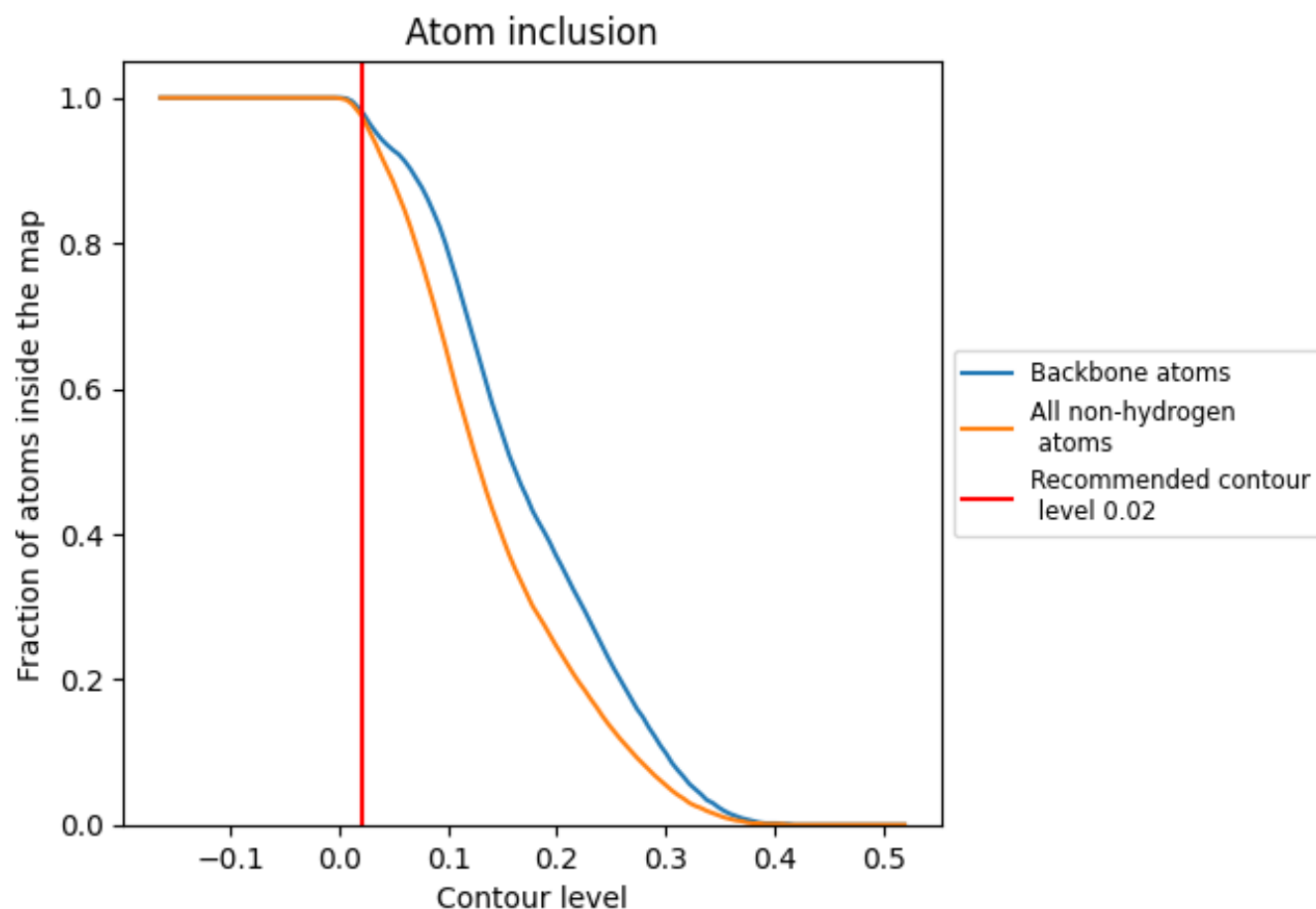
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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



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 (0.02).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 98% 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 (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9760	 0.4060
A	 0.9940	 0.4230
B	 0.9950	 0.4520
C	 0.9940	 0.4310
D	 0.7700	 0.0360
E	 0.6050	 0.0370
F	 1.0000	 0.3120
G	 1.0000	 0.1830
H	 1.0000	 0.2940
I	 1.0000	 0.2830
J	 1.0000	 0.3370
K	 1.0000	 0.3930
L	 0.9470	 0.1670
M	 1.0000	 0.3020
N	 1.0000	 0.2550
O	 1.0000	 0.3200
P	 1.0000	 0.3300
Q	 1.0000	 0.3060
R	 1.0000	 0.3140
S	 1.0000	 0.3390
T	 1.0000	 0.3810
U	 1.0000	 0.3580
V	 0.9470	 0.2790
W	 1.0000	 0.3430
X	 1.0000	 0.2970
Y	 1.0000	 0.3170
Z	 1.0000	 0.2230
a	 1.0000	 0.3390
b	 1.0000	 0.2300
c	 1.0000	 0.3020
d	 1.0000	 0.3180
e	 1.0000	 0.3910
f	 1.0000	 0.2970
g	 0.9740	 0.0980
h	 1.0000	 0.3610



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
i	 1.0000	 0.2230
j	 1.0000	 0.2330
k	 1.0000	 0.3070