



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 06:05 AM UTC

PDB ID : 8WQ0 / pdb_00008wq0
EMDB ID : EMD-37732
Title : Cryo-EM structure of WIV1 spike glycoprotein (the closed state)
Authors : Wang, X.; Qiao, S.
Deposited on : 2023-10-10
Resolution : 2.78 Å (reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

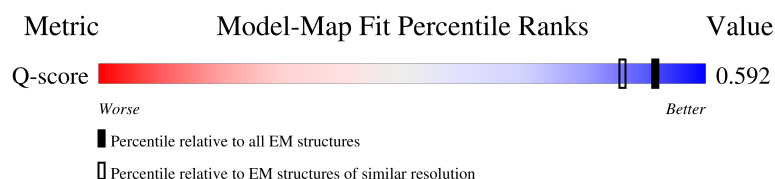
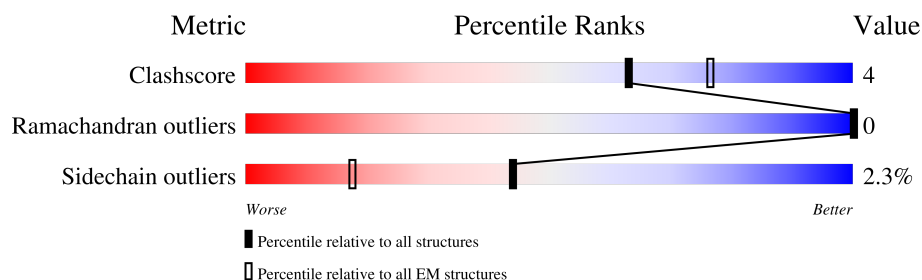
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13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

1 Overall quality at a glance

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

The reported resolution of this entry is 2.78 Å.

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



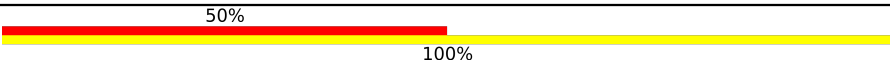
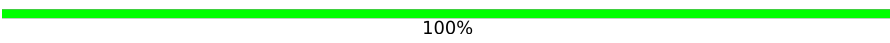
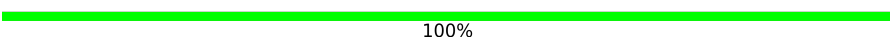
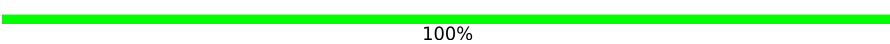
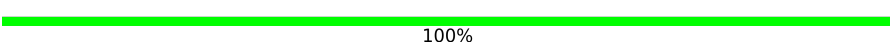
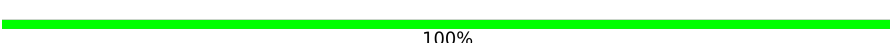
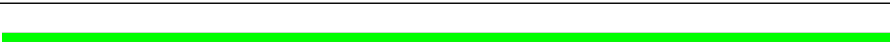
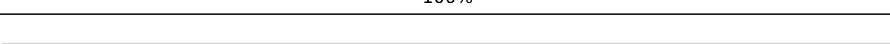
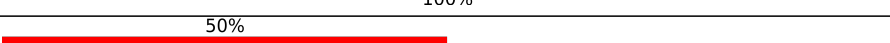
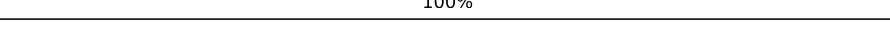
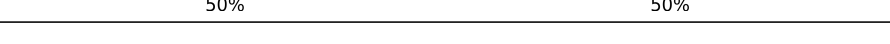
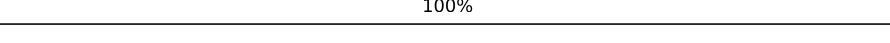
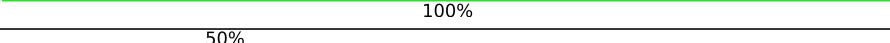
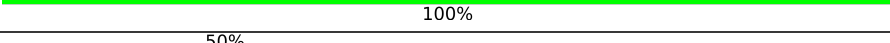
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10754 (2.28 - 3.28)

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	1271	
1	B	1271	
1	C	1271	
2	D	2	

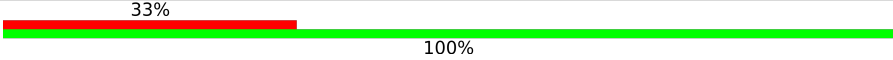
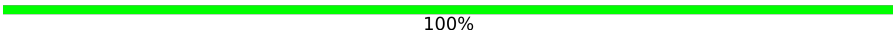
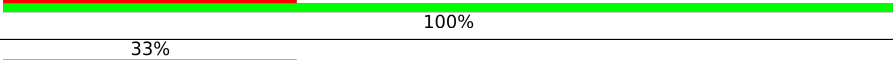
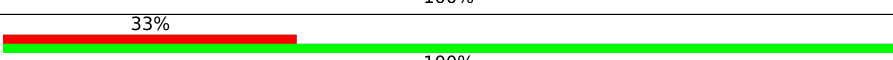
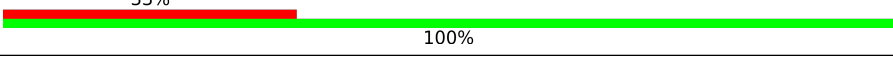
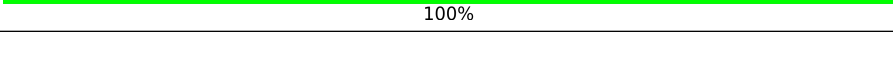
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Mol	Chain	Length	Quality of chain
2	F	2	
2	G	2	
2	H	2	
2	J	2	
2	K	2	
2	L	2	
2	M	2	
2	N	2	
2	O	2	
2	P	2	
2	R	2	
2	T	2	
2	U	2	
2	V	2	
2	W	2	
2	X	2	
2	Y	2	
2	Z	2	
2	b	2	
2	c	2	
2	d	2	
2	e	2	
2	g	2	
2	h	2	
2	i	2	

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Mol	Chain	Length	Quality of chain
3	E	3	
3	I	3	
3	Q	3	
3	S	3	
3	a	3	
3	f	3	
3	j	3	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 26750 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	1088	Total	C	N	O	S	0	0
			8485	5410	1414	1617	44		
1	B	1088	Total	C	N	O	S	0	0
			8485	5410	1414	1617	44		
1	C	1088	Total	C	N	O	S	0	0
			8485	5410	1414	1617	44		

There are 246 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	969	PRO	LYS	conflict	UNP U5WI05
A	970	PRO	VAL	conflict	UNP U5WI05
A	1192	GLY	-	expression tag	UNP U5WI05
A	1193	SER	-	expression tag	UNP U5WI05
A	1194	GLY	-	expression tag	UNP U5WI05
A	1195	TYR	-	expression tag	UNP U5WI05
A	1196	ILE	-	expression tag	UNP U5WI05
A	1197	PRO	-	expression tag	UNP U5WI05
A	1198	GLU	-	expression tag	UNP U5WI05
A	1199	ALA	-	expression tag	UNP U5WI05
A	1200	PRO	-	expression tag	UNP U5WI05
A	1201	ARG	-	expression tag	UNP U5WI05
A	1202	ASP	-	expression tag	UNP U5WI05
A	1203	GLY	-	expression tag	UNP U5WI05
A	1204	GLN	-	expression tag	UNP U5WI05
A	1205	ALA	-	expression tag	UNP U5WI05
A	1206	TYR	-	expression tag	UNP U5WI05
A	1207	VAL	-	expression tag	UNP U5WI05
A	1208	ARG	-	expression tag	UNP U5WI05
A	1209	LYS	-	expression tag	UNP U5WI05
A	1210	ASP	-	expression tag	UNP U5WI05
A	1211	GLY	-	expression tag	UNP U5WI05
A	1212	GLU	-	expression tag	UNP U5WI05
A	1213	TRP	-	expression tag	UNP U5WI05

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1214	VAL	-	expression tag	UNP U5WI05
A	1215	LEU	-	expression tag	UNP U5WI05
A	1216	LEU	-	expression tag	UNP U5WI05
A	1217	SER	-	expression tag	UNP U5WI05
A	1218	THR	-	expression tag	UNP U5WI05
A	1219	PHE	-	expression tag	UNP U5WI05
A	1220	LEU	-	expression tag	UNP U5WI05
A	1221	GLY	-	expression tag	UNP U5WI05
A	1222	ARG	-	expression tag	UNP U5WI05
A	1223	SER	-	expression tag	UNP U5WI05
A	1224	LEU	-	expression tag	UNP U5WI05
A	1225	GLU	-	expression tag	UNP U5WI05
A	1226	VAL	-	expression tag	UNP U5WI05
A	1227	LEU	-	expression tag	UNP U5WI05
A	1228	PHE	-	expression tag	UNP U5WI05
A	1229	GLN	-	expression tag	UNP U5WI05
A	1230	GLY	-	expression tag	UNP U5WI05
A	1231	PRO	-	expression tag	UNP U5WI05
A	1232	GLY	-	expression tag	UNP U5WI05
A	1233	HIS	-	expression tag	UNP U5WI05
A	1234	HIS	-	expression tag	UNP U5WI05
A	1235	HIS	-	expression tag	UNP U5WI05
A	1236	HIS	-	expression tag	UNP U5WI05
A	1237	HIS	-	expression tag	UNP U5WI05
A	1238	HIS	-	expression tag	UNP U5WI05
A	1239	HIS	-	expression tag	UNP U5WI05
A	1240	HIS	-	expression tag	UNP U5WI05
A	1241	SER	-	expression tag	UNP U5WI05
A	1242	ALA	-	expression tag	UNP U5WI05
A	1243	TRP	-	expression tag	UNP U5WI05
A	1244	SER	-	expression tag	UNP U5WI05
A	1245	HIS	-	expression tag	UNP U5WI05
A	1246	PRO	-	expression tag	UNP U5WI05
A	1247	GLN	-	expression tag	UNP U5WI05
A	1248	PHE	-	expression tag	UNP U5WI05
A	1249	GLU	-	expression tag	UNP U5WI05
A	1250	LYS	-	expression tag	UNP U5WI05
A	1251	GLY	-	expression tag	UNP U5WI05
A	1252	GLY	-	expression tag	UNP U5WI05
A	1253	GLY	-	expression tag	UNP U5WI05
A	1254	SER	-	expression tag	UNP U5WI05
A	1255	GLY	-	expression tag	UNP U5WI05

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1256	GLY	-	expression tag	UNP U5WI05
A	1257	GLY	-	expression tag	UNP U5WI05
A	1258	GLY	-	expression tag	UNP U5WI05
A	1259	SER	-	expression tag	UNP U5WI05
A	1260	GLY	-	expression tag	UNP U5WI05
A	1261	GLY	-	expression tag	UNP U5WI05
A	1262	SER	-	expression tag	UNP U5WI05
A	1263	ALA	-	expression tag	UNP U5WI05
A	1264	TRP	-	expression tag	UNP U5WI05
A	1265	SER	-	expression tag	UNP U5WI05
A	1266	HIS	-	expression tag	UNP U5WI05
A	1267	PRO	-	expression tag	UNP U5WI05
A	1268	GLN	-	expression tag	UNP U5WI05
A	1269	PHE	-	expression tag	UNP U5WI05
A	1270	GLU	-	expression tag	UNP U5WI05
A	1271	LYS	-	expression tag	UNP U5WI05
B	969	PRO	LYS	conflict	UNP U5WI05
B	970	PRO	VAL	conflict	UNP U5WI05
B	1192	GLY	-	expression tag	UNP U5WI05
B	1193	SER	-	expression tag	UNP U5WI05
B	1194	GLY	-	expression tag	UNP U5WI05
B	1195	TYR	-	expression tag	UNP U5WI05
B	1196	ILE	-	expression tag	UNP U5WI05
B	1197	PRO	-	expression tag	UNP U5WI05
B	1198	GLU	-	expression tag	UNP U5WI05
B	1199	ALA	-	expression tag	UNP U5WI05
B	1200	PRO	-	expression tag	UNP U5WI05
B	1201	ARG	-	expression tag	UNP U5WI05
B	1202	ASP	-	expression tag	UNP U5WI05
B	1203	GLY	-	expression tag	UNP U5WI05
B	1204	GLN	-	expression tag	UNP U5WI05
B	1205	ALA	-	expression tag	UNP U5WI05
B	1206	TYR	-	expression tag	UNP U5WI05
B	1207	VAL	-	expression tag	UNP U5WI05
B	1208	ARG	-	expression tag	UNP U5WI05
B	1209	LYS	-	expression tag	UNP U5WI05
B	1210	ASP	-	expression tag	UNP U5WI05
B	1211	GLY	-	expression tag	UNP U5WI05
B	1212	GLU	-	expression tag	UNP U5WI05
B	1213	TRP	-	expression tag	UNP U5WI05
B	1214	VAL	-	expression tag	UNP U5WI05
B	1215	LEU	-	expression tag	UNP U5WI05

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1216	LEU	-	expression tag	UNP U5WI05
B	1217	SER	-	expression tag	UNP U5WI05
B	1218	THR	-	expression tag	UNP U5WI05
B	1219	PHE	-	expression tag	UNP U5WI05
B	1220	LEU	-	expression tag	UNP U5WI05
B	1221	GLY	-	expression tag	UNP U5WI05
B	1222	ARG	-	expression tag	UNP U5WI05
B	1223	SER	-	expression tag	UNP U5WI05
B	1224	LEU	-	expression tag	UNP U5WI05
B	1225	GLU	-	expression tag	UNP U5WI05
B	1226	VAL	-	expression tag	UNP U5WI05
B	1227	LEU	-	expression tag	UNP U5WI05
B	1228	PHE	-	expression tag	UNP U5WI05
B	1229	GLN	-	expression tag	UNP U5WI05
B	1230	GLY	-	expression tag	UNP U5WI05
B	1231	PRO	-	expression tag	UNP U5WI05
B	1232	GLY	-	expression tag	UNP U5WI05
B	1233	HIS	-	expression tag	UNP U5WI05
B	1234	HIS	-	expression tag	UNP U5WI05
B	1235	HIS	-	expression tag	UNP U5WI05
B	1236	HIS	-	expression tag	UNP U5WI05
B	1237	HIS	-	expression tag	UNP U5WI05
B	1238	HIS	-	expression tag	UNP U5WI05
B	1239	HIS	-	expression tag	UNP U5WI05
B	1240	HIS	-	expression tag	UNP U5WI05
B	1241	SER	-	expression tag	UNP U5WI05
B	1242	ALA	-	expression tag	UNP U5WI05
B	1243	TRP	-	expression tag	UNP U5WI05
B	1244	SER	-	expression tag	UNP U5WI05
B	1245	HIS	-	expression tag	UNP U5WI05
B	1246	PRO	-	expression tag	UNP U5WI05
B	1247	GLN	-	expression tag	UNP U5WI05
B	1248	PHE	-	expression tag	UNP U5WI05
B	1249	GLU	-	expression tag	UNP U5WI05
B	1250	LYS	-	expression tag	UNP U5WI05
B	1251	GLY	-	expression tag	UNP U5WI05
B	1252	GLY	-	expression tag	UNP U5WI05
B	1253	GLY	-	expression tag	UNP U5WI05
B	1254	SER	-	expression tag	UNP U5WI05
B	1255	GLY	-	expression tag	UNP U5WI05
B	1256	GLY	-	expression tag	UNP U5WI05
B	1257	GLY	-	expression tag	UNP U5WI05

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1258	GLY	-	expression tag	UNP U5WI05
B	1259	SER	-	expression tag	UNP U5WI05
B	1260	GLY	-	expression tag	UNP U5WI05
B	1261	GLY	-	expression tag	UNP U5WI05
B	1262	SER	-	expression tag	UNP U5WI05
B	1263	ALA	-	expression tag	UNP U5WI05
B	1264	TRP	-	expression tag	UNP U5WI05
B	1265	SER	-	expression tag	UNP U5WI05
B	1266	HIS	-	expression tag	UNP U5WI05
B	1267	PRO	-	expression tag	UNP U5WI05
B	1268	GLN	-	expression tag	UNP U5WI05
B	1269	PHE	-	expression tag	UNP U5WI05
B	1270	GLU	-	expression tag	UNP U5WI05
B	1271	LYS	-	expression tag	UNP U5WI05
C	969	PRO	LYS	conflict	UNP U5WI05
C	970	PRO	VAL	conflict	UNP U5WI05
C	1192	GLY	-	expression tag	UNP U5WI05
C	1193	SER	-	expression tag	UNP U5WI05
C	1194	GLY	-	expression tag	UNP U5WI05
C	1195	TYR	-	expression tag	UNP U5WI05
C	1196	ILE	-	expression tag	UNP U5WI05
C	1197	PRO	-	expression tag	UNP U5WI05
C	1198	GLU	-	expression tag	UNP U5WI05
C	1199	ALA	-	expression tag	UNP U5WI05
C	1200	PRO	-	expression tag	UNP U5WI05
C	1201	ARG	-	expression tag	UNP U5WI05
C	1202	ASP	-	expression tag	UNP U5WI05
C	1203	GLY	-	expression tag	UNP U5WI05
C	1204	GLN	-	expression tag	UNP U5WI05
C	1205	ALA	-	expression tag	UNP U5WI05
C	1206	TYR	-	expression tag	UNP U5WI05
C	1207	VAL	-	expression tag	UNP U5WI05
C	1208	ARG	-	expression tag	UNP U5WI05
C	1209	LYS	-	expression tag	UNP U5WI05
C	1210	ASP	-	expression tag	UNP U5WI05
C	1211	GLY	-	expression tag	UNP U5WI05
C	1212	GLU	-	expression tag	UNP U5WI05
C	1213	TRP	-	expression tag	UNP U5WI05
C	1214	VAL	-	expression tag	UNP U5WI05
C	1215	LEU	-	expression tag	UNP U5WI05
C	1216	LEU	-	expression tag	UNP U5WI05
C	1217	SER	-	expression tag	UNP U5WI05

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1218	THR	-	expression tag	UNP U5WI05
C	1219	PHE	-	expression tag	UNP U5WI05
C	1220	LEU	-	expression tag	UNP U5WI05
C	1221	GLY	-	expression tag	UNP U5WI05
C	1222	ARG	-	expression tag	UNP U5WI05
C	1223	SER	-	expression tag	UNP U5WI05
C	1224	LEU	-	expression tag	UNP U5WI05
C	1225	GLU	-	expression tag	UNP U5WI05
C	1226	VAL	-	expression tag	UNP U5WI05
C	1227	LEU	-	expression tag	UNP U5WI05
C	1228	PHE	-	expression tag	UNP U5WI05
C	1229	GLN	-	expression tag	UNP U5WI05
C	1230	GLY	-	expression tag	UNP U5WI05
C	1231	PRO	-	expression tag	UNP U5WI05
C	1232	GLY	-	expression tag	UNP U5WI05
C	1233	HIS	-	expression tag	UNP U5WI05
C	1234	HIS	-	expression tag	UNP U5WI05
C	1235	HIS	-	expression tag	UNP U5WI05
C	1236	HIS	-	expression tag	UNP U5WI05
C	1237	HIS	-	expression tag	UNP U5WI05
C	1238	HIS	-	expression tag	UNP U5WI05
C	1239	HIS	-	expression tag	UNP U5WI05
C	1240	HIS	-	expression tag	UNP U5WI05
C	1241	SER	-	expression tag	UNP U5WI05
C	1242	ALA	-	expression tag	UNP U5WI05
C	1243	TRP	-	expression tag	UNP U5WI05
C	1244	SER	-	expression tag	UNP U5WI05
C	1245	HIS	-	expression tag	UNP U5WI05
C	1246	PRO	-	expression tag	UNP U5WI05
C	1247	GLN	-	expression tag	UNP U5WI05
C	1248	PHE	-	expression tag	UNP U5WI05
C	1249	GLU	-	expression tag	UNP U5WI05
C	1250	LYS	-	expression tag	UNP U5WI05
C	1251	GLY	-	expression tag	UNP U5WI05
C	1252	GLY	-	expression tag	UNP U5WI05
C	1253	GLY	-	expression tag	UNP U5WI05
C	1254	SER	-	expression tag	UNP U5WI05
C	1255	GLY	-	expression tag	UNP U5WI05
C	1256	GLY	-	expression tag	UNP U5WI05
C	1257	GLY	-	expression tag	UNP U5WI05
C	1258	GLY	-	expression tag	UNP U5WI05
C	1259	SER	-	expression tag	UNP U5WI05

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1260	GLY	-	expression tag	UNP U5WI05
C	1261	GLY	-	expression tag	UNP U5WI05
C	1262	SER	-	expression tag	UNP U5WI05
C	1263	ALA	-	expression tag	UNP U5WI05
C	1264	TRP	-	expression tag	UNP U5WI05
C	1265	SER	-	expression tag	UNP U5WI05
C	1266	HIS	-	expression tag	UNP U5WI05
C	1267	PRO	-	expression tag	UNP U5WI05
C	1268	GLN	-	expression tag	UNP U5WI05
C	1269	PHE	-	expression tag	UNP U5WI05
C	1270	GLU	-	expression tag	UNP U5WI05
C	1271	LYS	-	expression tag	UNP U5WI05

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



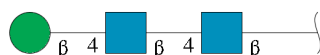
Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	2	Total	C	N	O	0	0
			28	16	2	10		
2	F	2	Total	C	N	O	0	0
			28	16	2	10		
2	G	2	Total	C	N	O	0	0
			28	16	2	10		
2	H	2	Total	C	N	O	0	0
			28	16	2	10		
2	J	2	Total	C	N	O	0	0
			28	16	2	10		
2	K	2	Total	C	N	O	0	0
			28	16	2	10		
2	L	2	Total	C	N	O	0	0
			28	16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

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Mol	Chain	Residues	Atoms				AltConf	Trace
2	P	2	Total	C	N	O	0	0
			28	16	2	10		
2	R	2	Total	C	N	O	0	0
			28	16	2	10		
2	T	2	Total	C	N	O	0	0
			28	16	2	10		
2	U	2	Total	C	N	O	0	0
			28	16	2	10		
2	V	2	Total	C	N	O	0	0
			28	16	2	10		
2	W	2	Total	C	N	O	0	0
			28	16	2	10		
2	X	2	Total	C	N	O	0	0
			28	16	2	10		
2	Y	2	Total	C	N	O	0	0
			28	16	2	10		
2	Z	2	Total	C	N	O	0	0
			28	16	2	10		
2	b	2	Total	C	N	O	0	0
			28	16	2	10		
2	c	2	Total	C	N	O	0	0
			28	16	2	10		
2	d	2	Total	C	N	O	0	0
			28	16	2	10		
2	e	2	Total	C	N	O	0	0
			28	16	2	10		
2	g	2	Total	C	N	O	0	0
			28	16	2	10		
2	h	2	Total	C	N	O	0	0
			28	16	2	10		
2	i	2	Total	C	N	O	0	0
			28	16	2	10		

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



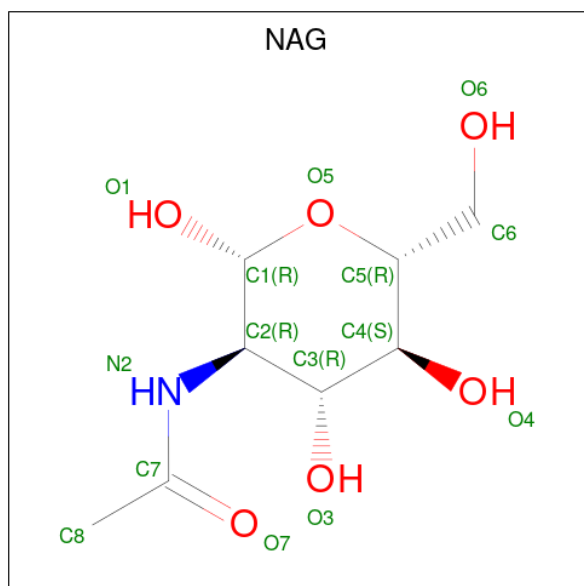
Mol	Chain	Residues	Atoms				AltConf	Trace
3	E	3	Total	C	N	O	0	0
			39	22	2	15		

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Mol	Chain	Residues	Atoms				AltConf	Trace
3	I	3	Total	C	N	O	0	0
			39	22	2	15		
3	Q	3	Total	C	N	O	0	0
			39	22	2	15		
3	S	3	Total	C	N	O	0	0
			39	22	2	15		
3	a	3	Total	C	N	O	0	0
			39	22	2	15		
3	f	3	Total	C	N	O	0	0
			39	22	2	15		
3	j	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

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



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

Chain L:  100%



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

Chain M:  100%



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

Chain N:  50% 50%



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

Chain O:  50% 100%



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

Chain P:  100%



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

Chain R:  50% 50%



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

Chain T:  100%



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

Chain U:  100%



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

Chain V:  100%



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

Chain W:  100%



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

Chain X:  100%



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

Chain Y:  50%  100%

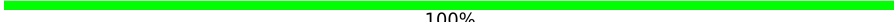


- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z:  50% 50%



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

Chain b:  100%



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

Chain c:  50% 50%



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

Chain d:  100%



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

Chain e:  50% 100%



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

Chain g:  50% 50% 50%



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

Chain h:  50% 50%



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

Chain i:  100%



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

Chain E:  100%



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

Chain I:  100%



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

Chain Q:  100%



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

Chain S:  100%



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

Chain a:  100%



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



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



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	307297	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.183	Depositor
Minimum map value	-0.090	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.0115	Depositor
Map size ($\text{\AA}$)	300.776, 300.776, 300.776	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.0742, 1.0742, 1.0742	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, 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	0.18	1/8689 (0.0%)	0.41	7/11828 (0.1%)
1	B	0.15	0/8689	0.33	0/11828
1	C	0.14	0/8689	0.34	2/11828 (0.0%)
All	All	0.16	1/26067 (0.0%)	0.36	9/35484 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	159	ASN	CA-C	-5.82	1.45	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	159	ASN	CA-C-N	-7.17	111.11	122.73
1	A	159	ASN	C-N-CA	-7.17	111.11	122.73
1	A	317	PHE	CA-C-N	-6.20	114.26	120.21
1	A	317	PHE	C-N-CA	-6.20	114.26	120.21
1	C	317	PHE	CA-C-N	-5.98	114.11	120.03
1	C	317	PHE	C-N-CA	-5.98	114.11	120.03
1	A	106	SER	N-CA-C	5.86	117.74	111.36
1	A	1122	ASP	CA-C-N	-5.06	114.45	119.56
1	A	1122	ASP	C-N-CA	-5.06	114.45	119.56

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	8485	0	8232	91	0
1	B	8485	0	8232	75	0
1	C	8485	0	8232	73	0
2	D	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	1	0
2	H	28	0	25	0	0
2	J	28	0	25	0	0
2	K	28	0	25	0	0
2	L	28	0	25	0	0
2	M	28	0	25	0	0
2	N	28	0	25	1	0
2	O	28	0	25	0	0
2	P	28	0	25	0	0
2	R	28	0	25	0	0
2	T	28	0	25	0	0
2	U	28	0	25	0	0
2	V	28	0	25	0	0
2	W	28	0	25	0	0
2	X	28	0	25	0	0
2	Y	28	0	25	0	0
2	Z	28	0	25	1	0
2	b	28	0	25	0	0
2	c	28	0	25	0	0
2	d	28	0	25	0	0
2	e	28	0	25	0	0
2	g	28	0	25	0	0
2	h	28	0	25	1	0
2	i	28	0	25	0	0
3	E	39	0	34	0	0
3	I	39	0	34	0	0
3	Q	39	0	34	0	0
3	S	39	0	34	0	0
3	a	39	0	34	0	0
3	f	39	0	34	0	0
3	j	39	0	34	0	0
4	A	98	0	91	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	112	0	104	1	0
4	C	84	0	78	1	0
All	All	26750	0	25857	223	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 4.

All (223) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:886:ALA:HB1	1:A:896:GLN:HG2	1.59	0.83
1:B:128:ALA:HB3	1:B:162:PHE:HB3	1.70	0.74
1:C:886:ALA:HB1	1:C:896:GLN:HG2	1.70	0.73
1:A:128:ALA:HB3	1:A:162:PHE:HB3	1.71	0.71
1:A:129:CYS:HA	1:A:160:CYS:HB2	1.71	0.71
1:A:951:SER:HB3	1:B:740:GLY:HA3	1.73	0.71
1:B:36:SER:HB3	1:B:65:SER:H	1.56	0.71
1:A:36:SER:HB3	1:A:65:SER:H	1.57	0.70
1:A:740:GLY:HA3	1:C:951:SER:HB3	1.74	0.69
1:C:36:SER:HB3	1:C:65:SER:H	1.56	0.69
1:A:1074:ARG:NH2	1:A:1101:ASP:O	2.28	0.66
1:C:787:GLN:NE2	1:C:918:GLN:OE1	2.28	0.66
1:B:787:GLN:NE2	1:B:918:GLN:OE1	2.30	0.64
1:B:886:ALA:HB1	1:B:896:GLN:HG2	1.78	0.64
1:A:880:PRO:HG2	1:A:883:MET:HE3	1.79	0.64
1:A:760:ASN:OD1	1:A:1002:ARG:NH1	2.30	0.63
1:A:627:ASN:N	1:A:627:ASN:HD22	1.96	0.63
1:A:464:ASP:OD1	1:A:465:GLY:N	2.34	0.60
1:A:383:VAL:HA	1:A:501:SER:O	2.01	0.60
1:C:550:GLN:O	1:C:564:ARG:NH2	2.34	0.60
1:A:627:ASN:N	1:A:627:ASN:ND2	2.50	0.59
1:A:167:LYS:H	1:A:167:LYS:HD3	1.67	0.59
1:B:27:THR:HB	1:B:76:ASN:HB3	1.83	0.59
1:A:520:LEU:HD21	1:A:572:LEU:HD11	1.84	0.59
1:A:141:VAL:HG21	1:A:173:LEU:HB3	1.86	0.58
1:A:787:GLN:NE2	1:A:918:GLN:OE1	2.37	0.58
1:A:550:GLN:O	1:A:564:ARG:NH2	2.34	0.58
1:B:98:VAL:HA	1:B:237:THR:HG22	1.85	0.58
1:B:818:LYS:NZ	1:B:831:ASP:OD2	2.36	0.58
1:A:1122:ASP:HB3	1:A:1125:GLN:HG2	1.85	0.58
1:B:550:GLN:O	1:B:564:ARG:NH2	2.31	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:741:SER:OG	1:C:743:CYS:SG	2.62	0.57
1:B:433:GLN:H	1:B:433:GLN:CD	2.13	0.56
1:A:1029:GLY:HA2	1:B:873:ALA:HB1	1.88	0.56
1:C:98:VAL:HA	1:C:237:THR:HG22	1.85	0.56
1:B:520:LEU:HD21	1:B:572:LEU:HD11	1.88	0.56
1:A:725:ILE:O	1:A:983:ARG:NH1	2.37	0.56
1:A:818:LYS:NZ	1:A:831:ASP:OD2	2.34	0.56
1:C:167:LYS:HD3	1:C:167:LYS:H	1.69	0.56
1:B:760:ASN:OD1	1:B:1002:ARG:NH1	2.38	0.55
1:C:791:ASP:HB3	1:C:794:LYS:HB2	1.87	0.55
1:C:880:PRO:HB2	1:C:883:MET:HG3	1.89	0.55
1:C:128:ALA:HB3	1:C:162:PHE:HB3	1.88	0.55
1:A:743:CYS:HB3	1:A:746:LEU:HD12	1.89	0.55
1:C:27:THR:HB	1:C:76:ASN:HB3	1.89	0.55
4:B:1304:NAG:H83	4:B:1304:NAG:H3	1.88	0.55
1:A:98:VAL:HA	1:A:237:THR:HG22	1.87	0.54
1:B:1029:GLY:HA2	1:C:873:ALA:HB1	1.90	0.54
4:C:1304:NAG:H83	4:C:1304:NAG:H3	1.88	0.54
4:A:1303:NAG:H83	4:A:1303:NAG:H3	1.88	0.54
1:C:643:VAL:HG23	1:C:678:TYR:HB3	1.89	0.54
1:C:1067:GLU:O	1:C:1069:LYS:HD3	2.07	0.54
1:B:1124:LEU:HG	1:C:1124:LEU:HD21	1.90	0.53
1:C:698:PRO:HA	1:C:1055:GLU:HA	1.90	0.53
1:B:951:SER:OG	1:C:740:GLY:HA3	2.07	0.53
1:A:117:ILE:HG23	1:A:126:ILE:HG12	1.91	0.53
1:A:873:ALA:HB1	1:C:1029:GLY:HA2	1.90	0.53
1:A:698:PRO:HA	1:A:1055:GLU:HA	1.89	0.53
1:B:117:ILE:HG23	1:B:126:ILE:HG12	1.91	0.53
1:A:27:THR:HB	1:A:76:ASN:HB3	1.90	0.53
1:A:72:THR:O	1:A:250:ALA:HA	2.09	0.53
1:A:968:ASP:OD2	1:C:373:THR:OG1	2.27	0.53
1:B:643:VAL:HG23	1:B:678:TYR:HB3	1.90	0.52
1:B:698:PRO:HA	1:B:1055:GLU:HA	1.90	0.52
1:A:869:TRP:HZ3	1:A:884:GLN:HG3	1.75	0.52
1:C:818:LYS:NZ	1:C:831:ASP:OD2	2.38	0.51
1:A:643:VAL:HG23	1:A:678:TYR:HB3	1.93	0.51
1:B:742:PHE:CE1	1:B:984:LEU:HD22	2.44	0.51
1:A:371:SER:OG	1:B:971:GLU:OE1	2.29	0.51
1:C:760:ASN:OD1	1:C:1002:ARG:NH1	2.41	0.51
1:A:197:VAL:HB	1:A:221:PHE:HB2	1.93	0.51
1:C:670:THR:HG23	1:C:672:GLN:H	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:725:ILE:O	1:B:983:ARG:NH1	2.44	0.50
1:B:318:PRO:HG2	1:B:320:ILE:HD11	1.94	0.50
1:B:371:SER:OG	1:C:968:ASP:OD1	2.25	0.50
1:C:742:PHE:CE1	1:C:984:LEU:HD22	2.46	0.50
1:A:968:ASP:OD1	1:C:371:SER:OG	2.26	0.50
1:C:103:VAL:HG22	1:C:116:ILE:HG12	1.94	0.50
1:A:670:THR:HG23	1:A:672:GLN:H	1.78	0.49
1:B:338:VAL:HG22	1:B:410:ASN:HB3	1.95	0.49
1:A:158:PHE:O	1:A:159:ASN:C	2.56	0.49
1:A:1112:VAL:HB	1:A:1115:ILE:HB	1.95	0.49
1:C:520:LEU:HD21	1:C:572:LEU:HD11	1.94	0.49
1:C:836:GLN:HB2	1:C:841:LEU:HD12	1.95	0.49
1:A:770:GLN:HG3	1:C:684:ALA:HB3	1.95	0.48
1:A:373:THR:OG1	1:B:968:ASP:OD2	2.30	0.48
1:A:210:LEU:HD12	1:A:211:PRO:HD2	1.95	0.48
1:A:364:THR:HB	1:A:423:ALA:HB3	1.94	0.48
1:A:306:PHE:HB3	1:A:580:GLY:HA3	1.96	0.48
1:A:954:GLY:O	1:A:978:ARG:NH1	2.47	0.48
1:B:167:LYS:H	1:B:167:LYS:HD3	1.79	0.48
1:C:117:ILE:HG23	1:C:126:ILE:HG12	1.96	0.48
1:C:1112:VAL:HB	1:C:1115:ILE:HB	1.96	0.48
1:C:127:ARG:HG2	1:C:129:CYS:SG	2.54	0.47
1:B:627:ASN:OD1	1:B:628:ASN:N	2.47	0.47
1:B:886:ALA:HB2	1:B:899:LEU:HD12	1.97	0.47
1:C:101:GLY:HA3	1:C:118:MET:HA	1.97	0.47
1:C:306:PHE:HB3	1:C:580:GLY:HA3	1.95	0.47
1:A:167:LYS:HD3	1:A:167:LYS:N	2.30	0.47
1:A:791:ASP:HB3	1:A:794:LYS:HB2	1.97	0.47
1:B:103:VAL:HG22	1:B:116:ILE:HG12	1.97	0.47
1:B:373:THR:OG1	1:C:968:ASP:OD2	2.32	0.47
1:C:187:VAL:HG13	1:C:258:LEU:HD11	1.97	0.47
1:C:794:LYS:NZ	1:C:803:ASP:OD2	2.44	0.47
1:B:210:LEU:HD12	1:B:211:PRO:HD2	1.97	0.47
1:C:325:PRO:HB2	1:C:328:GLU:HB2	1.98	0.46
1:B:836:GLN:HB2	1:B:841:LEU:HD12	1.96	0.46
1:B:1112:VAL:HB	1:B:1115:ILE:HB	1.97	0.46
1:C:565:ASP:HB3	1:C:568:THR:O	2.14	0.46
1:A:746:LEU:HD22	1:A:991:VAL:HG21	1.96	0.46
1:B:306:PHE:HB3	1:B:580:GLY:HA3	1.97	0.46
1:C:167:LYS:HD3	1:C:167:LYS:N	2.30	0.46
1:A:908:ASN:O	1:A:912:LYS:HE2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1086:PHE:CZ	2:G:1:NAG:H62	2.51	0.46
1:A:1093:PHE:CZ	1:A:1095:PRO:HG3	2.51	0.46
1:C:1086:PHE:CZ	2:h:1:NAG:H62	2.51	0.45
1:A:283:PRO:O	1:A:286:GLU:HG2	2.16	0.45
1:B:953:PHE:O	1:B:978:ARG:HB3	2.16	0.45
1:A:103:VAL:HG22	1:A:116:ILE:HG12	1.98	0.45
1:A:371:SER:OG	1:B:968:ASP:OD1	2.29	0.45
1:A:800:PHE:HE1	1:A:918:GLN:HG3	1.81	0.45
1:A:705:ILE:HG12	1:A:1048:VAL:HG22	1.98	0.45
1:B:141:VAL:HG21	1:B:173:LEU:HB3	1.99	0.45
1:C:210:LEU:HD12	1:C:211:PRO:HD2	1.99	0.45
1:B:604:CYS:HB3	1:B:629:VAL:HG12	1.99	0.45
1:B:791:ASP:HB3	1:B:794:LYS:HB2	1.98	0.45
1:A:753:ILE:O	1:A:757:GLN:HG2	2.17	0.45
1:B:134:CYS:SG	1:B:155:ASN:ND2	2.89	0.45
1:B:1093:PHE:CZ	1:B:1095:PRO:HG3	2.52	0.45
1:A:259:LYS:HB3	1:A:259:LYS:HE3	1.76	0.45
1:C:1093:PHE:CZ	1:C:1095:PRO:HG3	2.51	0.45
1:B:1069:LYS:HE2	1:B:1069:LYS:HB2	1.84	0.45
1:A:21:LEU:HB2	1:A:138:PHE:HZ	1.82	0.44
1:A:134:CYS:SG	1:A:155:ASN:ND2	2.91	0.44
1:B:383:VAL:HA	1:B:501:SER:O	2.17	0.44
1:C:42:TYR:OH	1:C:59:HIS:O	2.24	0.44
1:A:325:PRO:HB2	1:A:328:GLU:HB2	1.99	0.44
1:B:954:GLY:O	1:B:978:ARG:NH1	2.51	0.44
1:C:141:VAL:HG21	1:C:173:LEU:HB3	2.00	0.44
1:B:101:GLY:HA3	1:B:118:MET:HA	2.00	0.44
1:A:470:PRO:HG3	1:A:477:TRP:CZ2	2.52	0.44
1:C:21:LEU:HB2	1:C:138:PHE:HZ	1.82	0.44
1:A:42:TYR:OH	1:A:59:HIS:O	2.23	0.44
1:A:868:GLY:HA2	1:A:884:GLN:OE1	2.18	0.44
1:A:1067:GLU:O	1:A:1069:LYS:HD3	2.17	0.44
1:B:470:PRO:HG3	1:B:477:TRP:CZ2	2.53	0.44
1:A:338:VAL:HG22	1:A:410:ASN:HB3	1.99	0.44
1:A:565:ASP:HB3	1:A:568:THR:O	2.18	0.44
1:B:325:PRO:HB2	1:B:328:GLU:HB2	2.00	0.43
1:B:338:VAL:HA	1:B:388:PHE:HB2	1.99	0.43
1:B:565:ASP:HB3	1:B:568:THR:O	2.18	0.43
1:C:470:PRO:HG3	1:C:477:TRP:CZ2	2.53	0.43
1:C:693:ASN:N	1:C:693:ASN:OD1	2.51	0.43
1:C:828:ASN:OD1	1:C:829:ALA:N	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:VAL:HA	1:B:128:ALA:HA	2.00	0.43
1:A:127:ARG:HG2	1:A:129:CYS:SG	2.58	0.43
1:A:1031:HIS:HA	1:A:1049:THR:HG22	1.99	0.43
1:B:259:LYS:HB3	1:B:259:LYS:HE3	1.76	0.43
1:B:308:VAL:HB	1:B:577:CYS:HB3	2.00	0.43
1:B:684:ALA:HB3	1:C:770:GLN:HG2	2.01	0.43
1:B:1086:PHE:HE1	1:B:1097:ILE:HD13	1.83	0.43
1:A:312:LYS:HD3	1:A:521:ILE:HD13	2.00	0.43
1:A:67:VAL:CG1	1:A:254:PHE:HB3	2.49	0.43
1:B:21:LEU:HB2	1:B:138:PHE:HZ	1.84	0.43
1:A:153:ILE:HD12	1:A:235:LEU:HD13	2.01	0.43
1:A:746:LEU:HD21	1:A:988:GLN:OE1	2.19	0.43
1:B:732:CYS:HB2	1:B:960:LEU:HD21	2.01	0.43
1:B:187:VAL:HG13	1:B:258:LEU:HD11	2.01	0.42
1:C:338:VAL:HA	1:C:388:PHE:HB2	2.00	0.42
1:A:196:HIS:CD2	1:A:222:LYS:HG3	2.54	0.42
1:C:945:LEU:HD12	1:C:945:LEU:HA	1.90	0.42
1:C:746:LEU:HD22	1:C:991:VAL:HG21	2.01	0.42
1:C:156:ASN:HD22	2:Z:1:NAG:H81	1.82	0.42
1:B:880:PRO:HB2	1:B:883:MET:HG3	2.01	0.42
1:A:186:PHE:HA	1:A:198:TYR:O	2.19	0.42
1:A:386:ASP:O	1:A:498:VAL:HA	2.18	0.42
1:A:707:THR:HG23	1:A:917:ILE:HD12	2.01	0.42
1:A:1081:ASN:OD1	1:A:1082:GLY:N	2.53	0.42
1:A:1124:LEU:HG	1:B:1124:LEU:HD21	2.00	0.42
1:B:710:MET:HE2	1:B:710:MET:HB3	1.84	0.42
1:A:945:LEU:HD12	1:A:945:LEU:HA	1.92	0.42
1:B:155:ASN:N	1:B:155:ASN:HD22	2.17	0.42
1:B:632:THR:OG1	1:B:635:GLY:N	2.42	0.42
1:C:883:MET:HE2	1:C:883:MET:HB3	1.80	0.42
1:A:836:GLN:HB2	1:A:841:LEU:HD12	2.01	0.42
1:A:101:GLY:HA3	1:A:118:MET:HA	2.02	0.42
1:A:953:PHE:CD2	1:A:982:GLY:HA3	2.55	0.42
1:C:259:LYS:HB3	1:C:259:LYS:HE3	1.76	0.42
1:C:627:ASN:OD1	1:C:628:ASN:N	2.52	0.42
1:C:1081:ASN:OD1	1:C:1082:GLY:N	2.52	0.42
1:A:308:VAL:HB	1:A:577:CYS:HB3	2.01	0.41
1:A:732:CYS:HB2	1:A:960:LEU:HD21	2.02	0.41
1:B:287:LEU:HD21	1:B:296:ILE:HG21	2.01	0.41
1:C:312:LYS:HD3	1:C:521:ILE:HD13	2.02	0.41
1:A:693:ASN:OD1	1:A:693:ASN:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:732:CYS:HB2	1:C:960:LEU:HD21	2.01	0.41
1:B:817:MET:HE2	1:B:817:MET:HB3	1.99	0.41
1:C:888:ARG:NH1	1:C:1032:LEU:O	2.53	0.41
1:A:336:PRO:HG3	1:A:342:GLU:HB2	2.02	0.41
1:A:904:LYS:HB3	1:A:904:LYS:HE2	1.88	0.41
1:B:129:CYS:HB2	1:B:131:PHE:CZ	2.56	0.41
1:C:954:GLY:O	1:C:978:ARG:NH1	2.53	0.41
1:A:794:LYS:HB2	1:A:794:LYS:HE3	1.86	0.41
1:B:49:ARG:O	1:B:271:GLY:HA2	2.20	0.41
1:C:904:LYS:HB3	1:C:904:LYS:HE2	1.89	0.41
1:C:918:GLN:O	1:C:922:THR:OG1	2.38	0.41
1:B:283:PRO:O	1:B:286:GLU:HG2	2.21	0.41
1:B:693:ASN:OD1	1:B:693:ASN:N	2.53	0.41
1:C:953:PHE:O	1:C:978:ARG:HB3	2.21	0.41
1:C:969:PRO:N	1:C:970:PRO:HD2	2.36	0.41
1:B:42:TYR:OH	1:B:59:HIS:O	2.23	0.41
1:C:189:ARG:NH2	1:C:191:LYS:HD3	2.36	0.40
1:A:100:ARG:HH22	1:A:173:LEU:HD11	1.86	0.40
1:A:567:LYS:HD3	2:N:1:NAG:H5	2.03	0.40
1:A:685:ASP:OD1	1:B:773:LYS:NZ	2.40	0.40
1:B:945:LEU:HD12	1:B:945:LEU:HA	1.90	0.40
1:C:318:PRO:HD2	1:C:512:CYS:SG	2.61	0.40
1:C:450:ARG:HB2	1:C:453:GLU:CD	2.47	0.40
1:B:223:LEU:HA	1:B:224:PRO:HD3	1.98	0.40
1:B:403:THR:HG22	1:C:372:ALA:HB1	2.04	0.40
1:B:826:ASP:OD2	1:B:829:ALA:HB3	2.22	0.40
1:C:386:ASP:O	1:C:498:VAL:HA	2.22	0.40
1:C:703:ILE:HG13	1:C:906:ILE:HG23	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1084/1271 (85%)	1054 (97%)	30 (3%)	0	100	100
1	B	1084/1271 (85%)	1056 (97%)	28 (3%)	0	100	100
1	C	1084/1271 (85%)	1060 (98%)	24 (2%)	0	100	100
All	All	3252/3813 (85%)	3170 (98%)	82 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	946/1100 (86%)	919 (97%)	27 (3%)	37	70
1	B	946/1100 (86%)	927 (98%)	19 (2%)	48	77
1	C	946/1100 (86%)	928 (98%)	18 (2%)	50	78
All	All	2838/3300 (86%)	2774 (98%)	64 (2%)	44	75

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	THR
1	A	55	LEU
1	A	93	THR
1	A	121	SER
1	A	141	VAL
1	A	160	CYS
1	A	236	LEU
1	A	291	VAL
1	A	358	ASN
1	A	383	VAL
1	A	501	SER
1	A	562	SER
1	A	595	VAL
1	A	627	ASN
1	A	672	GLN

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Mol	Chain	Res	Type
1	A	721	CYS
1	A	723	MET
1	A	776	THR
1	A	904	LYS
1	A	935	VAL
1	A	988	GLN
1	A	1026	CYS
1	A	1102	ASN
1	A	1105	VAL
1	A	1106	SER
1	A	1119	THR
1	A	1124	LEU
1	B	55	LEU
1	B	121	SER
1	B	130	ASN
1	B	141	VAL
1	B	291	VAL
1	B	358	ASN
1	B	383	VAL
1	B	501	SER
1	B	562	SER
1	B	595	VAL
1	B	604	CYS
1	B	776	THR
1	B	904	LYS
1	B	926	THR
1	B	935	VAL
1	B	1026	CYS
1	B	1077	VAL
1	B	1105	VAL
1	B	1119	THR
1	C	27	THR
1	C	55	LEU
1	C	113	GLN
1	C	121	SER
1	C	141	VAL
1	C	291	VAL
1	C	323	LEU
1	C	358	ASN
1	C	501	SER
1	C	562	SER
1	C	595	VAL

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Mol	Chain	Res	Type
1	C	776	THR
1	C	904	LYS
1	C	926	THR
1	C	935	VAL
1	C	1102	ASN
1	C	1105	VAL
1	C	1119	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	HIS
1	A	155	ASN
1	A	231	ASN
1	A	628	ASN
1	A	642	HIS
1	A	722	ASN
1	A	994	GLN
1	A	1089	GLN
1	B	113	GLN
1	B	155	ASN
1	B	628	ASN
1	B	902	ASN
1	C	113	GLN
1	C	382	ASN
1	C	628	ASN
1	C	642	HIS
1	C	672	GLN
1	C	1089	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

73 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	D	1	2,1	14,14,15	0.29	0	17,19,21	0.53	0
2	NAG	D	2	2	14,14,15	0.27	0	17,19,21	0.59	0
3	NAG	E	1	3,1	14,14,15	0.26	0	17,19,21	0.54	0
3	NAG	E	2	3	14,14,15	0.28	0	17,19,21	0.47	0
3	BMA	E	3	3	11,11,12	0.21	0	15,15,17	0.54	0
2	NAG	F	1	2,1	14,14,15	0.39	0	17,19,21	1.21	2 (11%)
2	NAG	F	2	2	14,14,15	0.42	0	17,19,21	1.63	5 (29%)
2	NAG	G	1	2,1	14,14,15	0.97	1 (7%)	17,19,21	1.04	1 (5%)
2	NAG	G	2	2	14,14,15	0.23	0	17,19,21	0.52	0
2	NAG	H	1	2,1	14,14,15	0.22	0	17,19,21	0.42	0
2	NAG	H	2	2	14,14,15	0.23	0	17,19,21	0.42	0
3	NAG	I	1	3,1	14,14,15	0.30	0	17,19,21	0.64	0
3	NAG	I	2	3	14,14,15	0.28	0	17,19,21	0.54	0
3	BMA	I	3	3	11,11,12	0.21	0	15,15,17	0.52	0
2	NAG	J	1	2,1	14,14,15	0.28	0	17,19,21	0.67	0
2	NAG	J	2	2	14,14,15	0.28	0	17,19,21	0.55	0
2	NAG	K	1	2,1	14,14,15	0.31	0	17,19,21	0.71	0
2	NAG	K	2	2	14,14,15	0.28	0	17,19,21	0.68	0
2	NAG	L	1	2,1	14,14,15	0.30	0	17,19,21	0.52	0
2	NAG	L	2	2	14,14,15	0.28	0	17,19,21	0.60	0
2	NAG	M	1	2,1	14,14,15	0.29	0	17,19,21	0.54	0
2	NAG	M	2	2	14,14,15	0.29	0	17,19,21	0.62	0
2	NAG	N	1	2,1	14,14,15	0.29	0	17,19,21	0.43	0
2	NAG	N	2	2	14,14,15	0.25	0	17,19,21	0.53	0
2	NAG	O	1	2,1	14,14,15	0.28	0	17,19,21	0.64	0
2	NAG	O	2	2	14,14,15	0.28	0	17,19,21	0.63	0
2	NAG	P	1	2,1	14,14,15	0.28	0	17,19,21	0.49	0
2	NAG	P	2	2	14,14,15	0.28	0	17,19,21	0.60	0
3	NAG	Q	1	3,1	14,14,15	0.27	0	17,19,21	0.56	0
3	NAG	Q	2	3	14,14,15	0.27	0	17,19,21	0.58	0
3	BMA	Q	3	3	11,11,12	0.22	0	15,15,17	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	R	1	2,1	14,14,15	0.28	0	17,19,21	0.86	1 (5%)
2	NAG	R	2	2	14,14,15	0.27	0	17,19,21	0.63	0
3	NAG	S	1	3,1	14,14,15	0.23	0	17,19,21	0.40	0
3	NAG	S	2	3	14,14,15	0.24	0	17,19,21	0.45	0
3	BMA	S	3	3	11,11,12	0.56	0	15,15,17	0.67	0
2	NAG	T	1	2,1	14,14,15	0.24	0	17,19,21	0.43	0
2	NAG	T	2	2	14,14,15	0.25	0	17,19,21	0.41	0
2	NAG	U	1	2,1	14,14,15	0.31	0	17,19,21	0.65	0
2	NAG	U	2	2	14,14,15	0.26	0	17,19,21	0.60	0
2	NAG	V	1	2,1	14,14,15	0.31	0	17,19,21	0.73	0
2	NAG	V	2	2	14,14,15	0.29	0	17,19,21	0.64	0
2	NAG	W	1	2,1	14,14,15	0.29	0	17,19,21	0.50	0
2	NAG	W	2	2	14,14,15	0.27	0	17,19,21	0.59	0
2	NAG	X	1	2,1	14,14,15	0.29	0	17,19,21	0.49	0
2	NAG	X	2	2	14,14,15	0.28	0	17,19,21	0.62	0
2	NAG	Y	1	2,1	14,14,15	0.29	0	17,19,21	0.64	0
2	NAG	Y	2	2	14,14,15	0.28	0	17,19,21	0.61	0
2	NAG	Z	1	2,1	14,14,15	0.26	0	17,19,21	0.49	0
2	NAG	Z	2	2	14,14,15	0.28	0	17,19,21	0.60	0
3	NAG	a	1	3,1	14,14,15	0.26	0	17,19,21	0.57	0
3	NAG	a	2	3	14,14,15	0.28	0	17,19,21	0.51	0
3	BMA	a	3	3	11,11,12	0.21	0	15,15,17	0.54	0
2	NAG	b	1	2,1	14,14,15	0.25	0	17,19,21	0.46	0
2	NAG	b	2	2	14,14,15	0.20	0	17,19,21	0.41	0
2	NAG	c	1	2,1	14,14,15	0.24	0	17,19,21	0.98	1 (5%)
2	NAG	c	2	2	14,14,15	0.29	0	17,19,21	0.63	0
2	NAG	d	1	2,1	14,14,15	0.29	0	17,19,21	0.47	0
2	NAG	d	2	2	14,14,15	0.27	0	17,19,21	0.56	0
2	NAG	e	1	2,1	14,14,15	0.28	0	17,19,21	0.59	0
2	NAG	e	2	2	14,14,15	0.28	0	17,19,21	0.59	0
3	NAG	f	1	3,1	14,14,15	0.23	0	17,19,21	0.40	0
3	NAG	f	2	3	14,14,15	0.24	0	17,19,21	0.44	0
3	BMA	f	3	3	11,11,12	0.59	0	15,15,17	0.69	0
2	NAG	g	1	2,1	14,14,15	0.51	0	17,19,21	0.70	1 (5%)
2	NAG	g	2	2	14,14,15	0.28	0	17,19,21	0.42	0
2	NAG	h	1	2,1	14,14,15	0.99	1 (7%)	17,19,21	1.04	1 (5%)
2	NAG	h	2	2	14,14,15	0.26	0	17,19,21	0.54	0
2	NAG	i	1	2,1	14,14,15	0.23	0	17,19,21	0.43	0
2	NAG	i	2	2	14,14,15	0.23	0	17,19,21	0.42	0
3	NAG	j	1	3,1	14,14,15	0.29	0	17,19,21	0.45	0
3	NAG	j	2	3	14,14,15	0.25	0	17,19,21	0.41	0
3	BMA	j	3	3	11,11,12	0.58	0	15,15,17	0.68	0

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

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

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	h	1	NAG	O5-C1	-3.35	1.38	1.43
2	G	1	NAG	O5-C1	-3.26	1.38	1.43

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	NAG	C2-N2-C7	3.55	127.66	122.90
2	F	2	NAG	O5-C1-C2	-2.92	106.78	111.29
2	F	1	NAG	O4-C4-C5	2.90	116.46	109.32
2	F	1	NAG	C1-O5-C5	2.79	115.92	112.19
2	G	1	NAG	C3-C4-C5	2.75	115.22	110.23
2	h	1	NAG	C3-C4-C5	2.74	115.20	110.23
2	c	1	NAG	C1-O5-C5	2.61	115.69	112.19
2	R	1	NAG	C4-C3-C2	-2.60	107.20	111.02
2	F	2	NAG	C1-O5-C5	2.38	115.38	112.19
2	F	2	NAG	C1-C2-N2	2.31	114.08	110.43
2	g	1	NAG	C1-O5-C5	2.10	115.00	112.19
2	F	2	NAG	O5-C5-C6	2.00	111.56	107.66

There are no chirality outliers.

All (81) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	2	NAG	C1-C2-N2-C7
2	M	2	NAG	C8-C7-N2-C2
2	M	2	NAG	O7-C7-N2-C2
2	P	1	NAG	C8-C7-N2-C2
2	P	1	NAG	O7-C7-N2-C2
2	Z	1	NAG	C8-C7-N2-C2
2	c	1	NAG	C1-C2-N2-C7
2	c	1	NAG	O7-C7-N2-C2
3	E	1	NAG	C8-C7-N2-C2
3	E	1	NAG	O7-C7-N2-C2
3	E	2	NAG	C8-C7-N2-C2
3	E	2	NAG	O7-C7-N2-C2
3	Q	1	NAG	C8-C7-N2-C2
3	Q	1	NAG	O7-C7-N2-C2
3	a	1	NAG	C8-C7-N2-C2
3	a	1	NAG	O7-C7-N2-C2
2	N	2	NAG	C8-C7-N2-C2
2	N	2	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	Z	1	NAG	O7-C7-N2-C2
2	c	1	NAG	C8-C7-N2-C2
2	d	2	NAG	C8-C7-N2-C2
2	d	2	NAG	O7-C7-N2-C2
2	R	2	NAG	C8-C7-N2-C2
2	R	2	NAG	O7-C7-N2-C2
2	i	1	NAG	O5-C5-C6-O6
2	H	1	NAG	O5-C5-C6-O6
2	U	1	NAG	C8-C7-N2-C2
2	U	1	NAG	O7-C7-N2-C2
2	T	1	NAG	O5-C5-C6-O6
2	H	1	NAG	C4-C5-C6-O6
2	T	1	NAG	C4-C5-C6-O6
2	K	1	NAG	C8-C7-N2-C2
2	N	1	NAG	C8-C7-N2-C2
2	V	1	NAG	C8-C7-N2-C2
2	d	1	NAG	C8-C7-N2-C2
3	I	1	NAG	C8-C7-N2-C2
2	H	2	NAG	C4-C5-C6-O6
3	j	2	NAG	C4-C5-C6-O6
2	i	1	NAG	C4-C5-C6-O6
2	D	1	NAG	C8-C7-N2-C2
2	K	1	NAG	O7-C7-N2-C2
2	N	1	NAG	O7-C7-N2-C2
2	O	1	NAG	C8-C7-N2-C2
2	R	1	NAG	C8-C7-N2-C2
2	V	1	NAG	O7-C7-N2-C2
2	Y	1	NAG	C8-C7-N2-C2
2	d	1	NAG	O7-C7-N2-C2
3	I	1	NAG	O7-C7-N2-C2
2	O	1	NAG	O7-C7-N2-C2
2	b	2	NAG	C4-C5-C6-O6
2	i	2	NAG	C4-C5-C6-O6
2	b	2	NAG	O5-C5-C6-O6
2	b	1	NAG	O5-C5-C6-O6
2	e	1	NAG	C8-C7-N2-C2
2	b	1	NAG	C4-C5-C6-O6
2	D	1	NAG	O7-C7-N2-C2
2	R	1	NAG	O7-C7-N2-C2
2	Y	1	NAG	O7-C7-N2-C2
3	j	2	NAG	O5-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6

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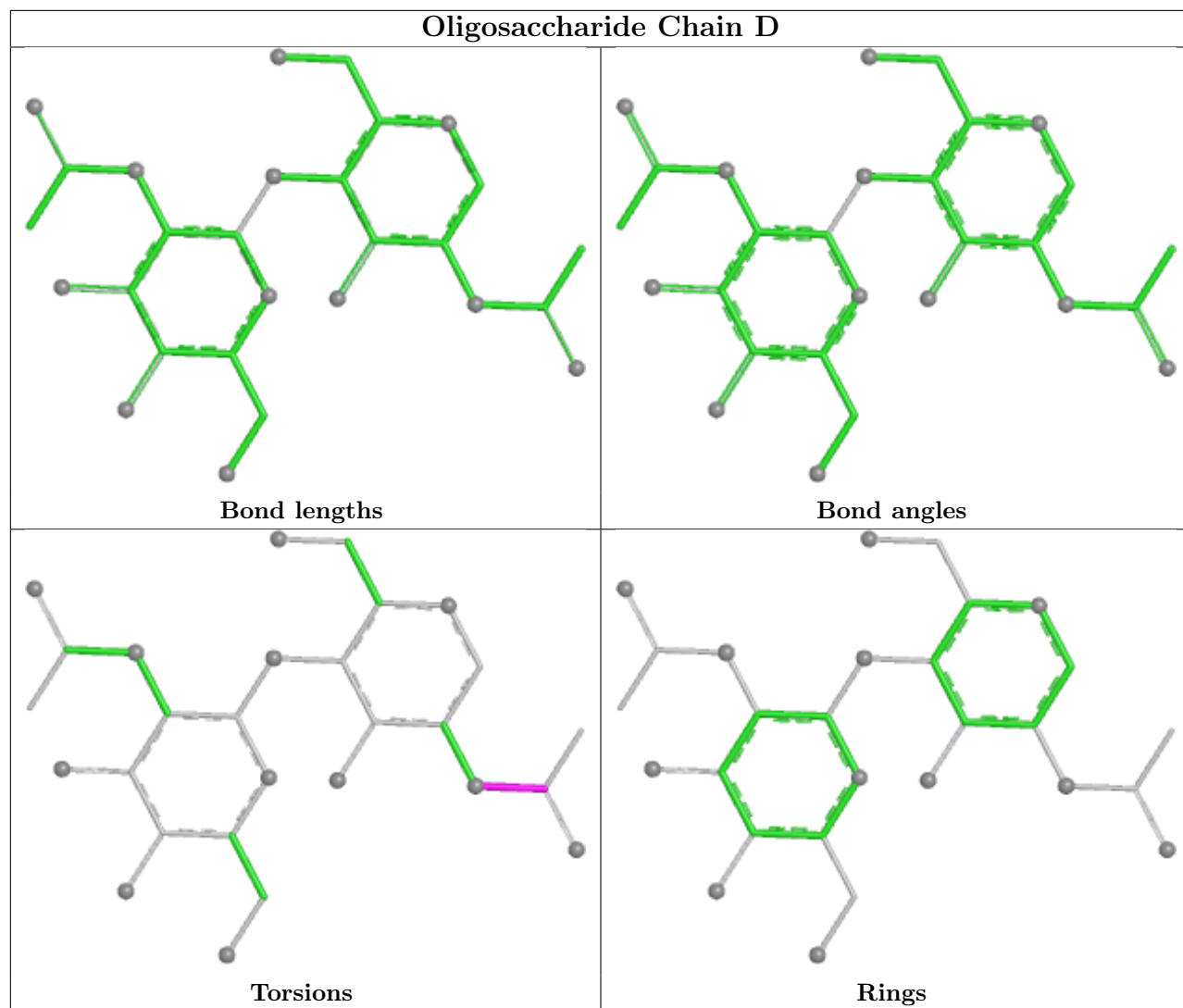
Mol	Chain	Res	Type	Atoms
2	g	1	NAG	O5-C5-C6-O6
2	g	1	NAG	C4-C5-C6-O6
2	e	1	NAG	O7-C7-N2-C2
2	T	2	NAG	C4-C5-C6-O6
3	f	2	NAG	C4-C5-C6-O6
3	f	2	NAG	O5-C5-C6-O6
3	S	2	NAG	C4-C5-C6-O6
2	i	2	NAG	O5-C5-C6-O6
3	S	2	NAG	O5-C5-C6-O6
2	M	2	NAG	O5-C5-C6-O6
2	W	2	NAG	C8-C7-N2-C2
2	T	2	NAG	O5-C5-C6-O6
2	W	1	NAG	C8-C7-N2-C2
2	W	2	NAG	O7-C7-N2-C2
2	W	1	NAG	O7-C7-N2-C2
2	M	1	NAG	C1-C2-N2-C7
2	N	1	NAG	C1-C2-N2-C7
2	W	2	NAG	C1-C2-N2-C7
2	L	1	NAG	C8-C7-N2-C2
3	f	1	NAG	C4-C5-C6-O6
2	L	1	NAG	O7-C7-N2-C2

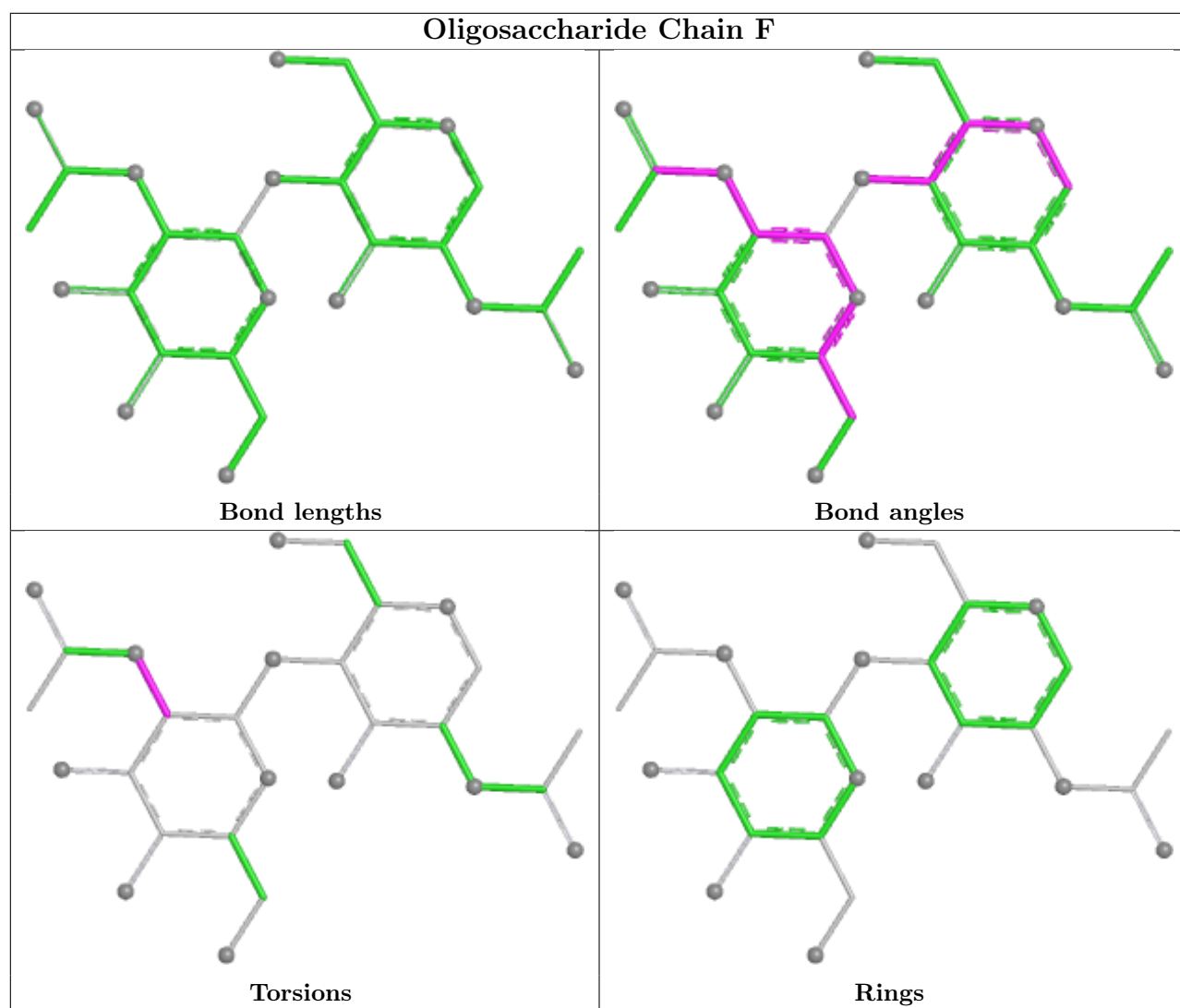
There are no ring outliers.

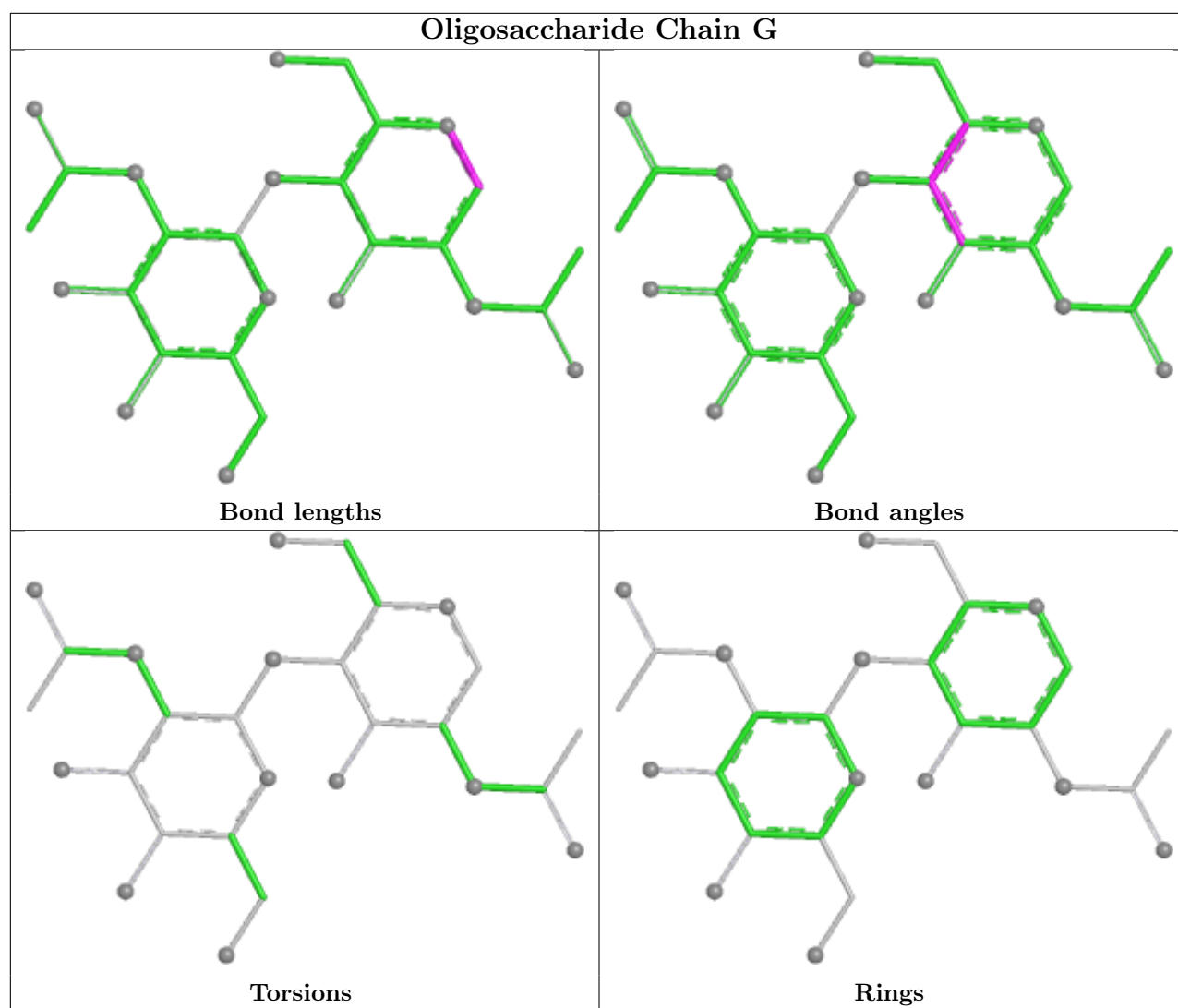
4 monomers are involved in 4 short contacts:

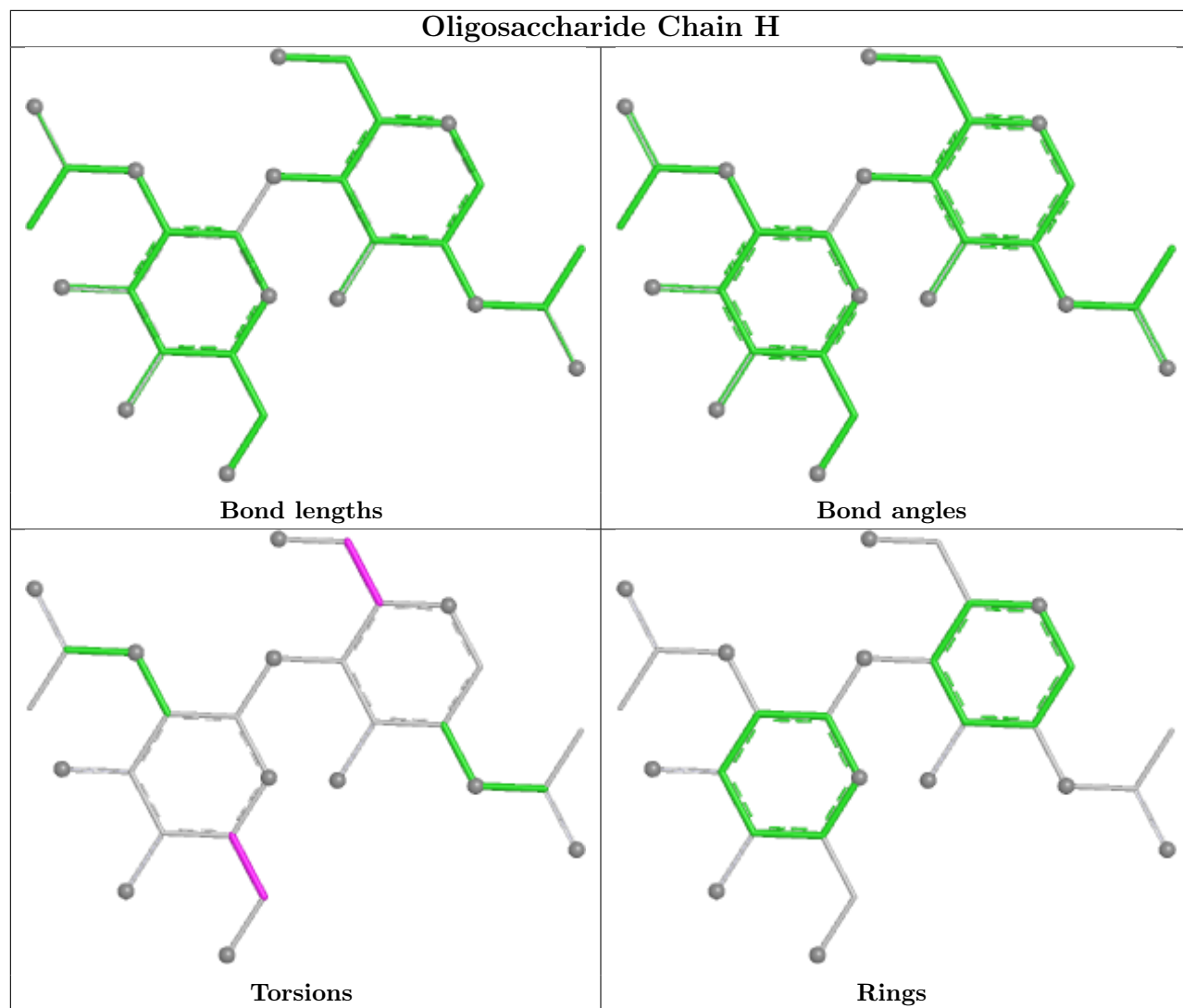
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	Z	1	NAG	1	0
2	h	1	NAG	1	0
2	G	1	NAG	1	0
2	N	1	NAG	1	0

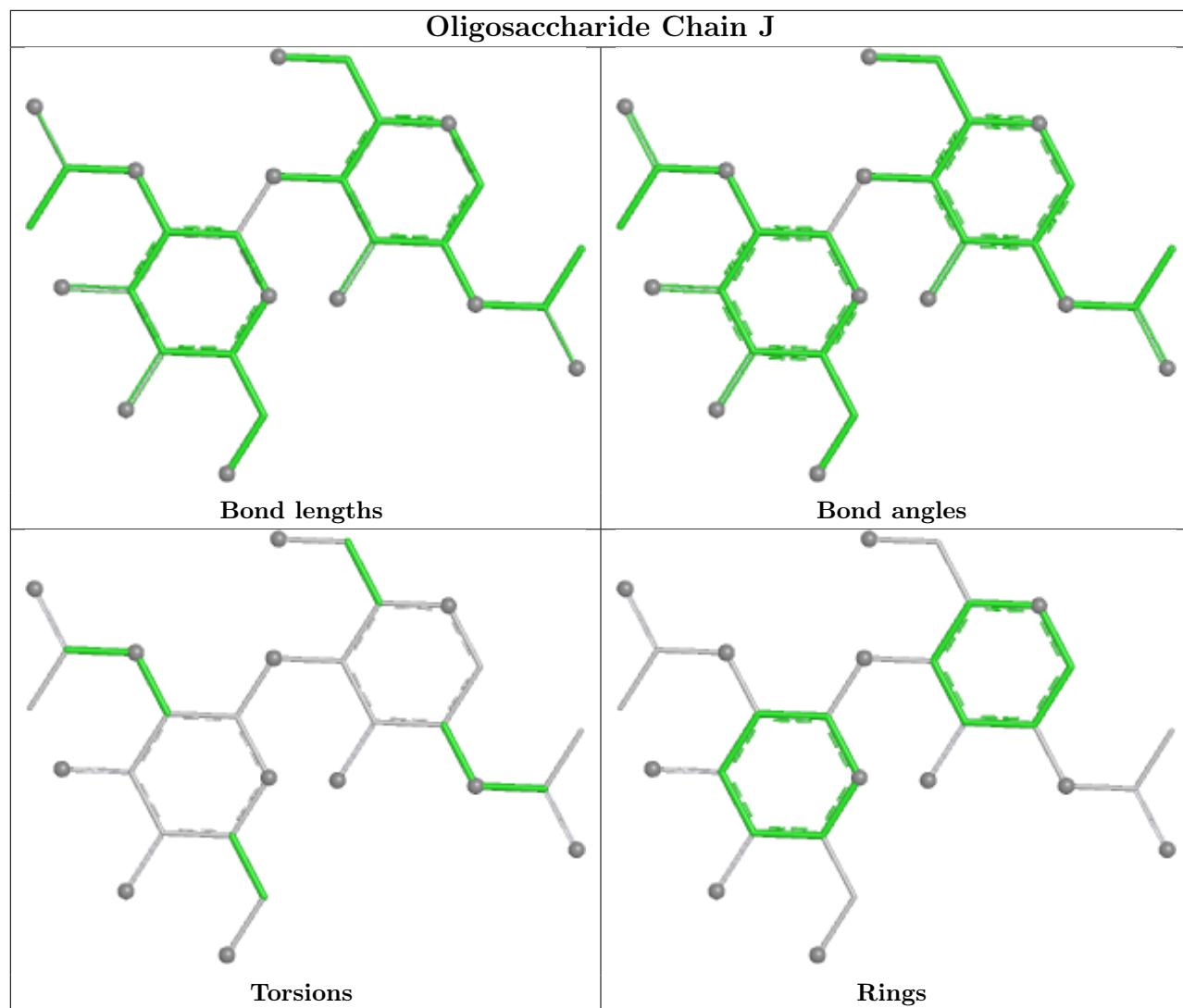
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

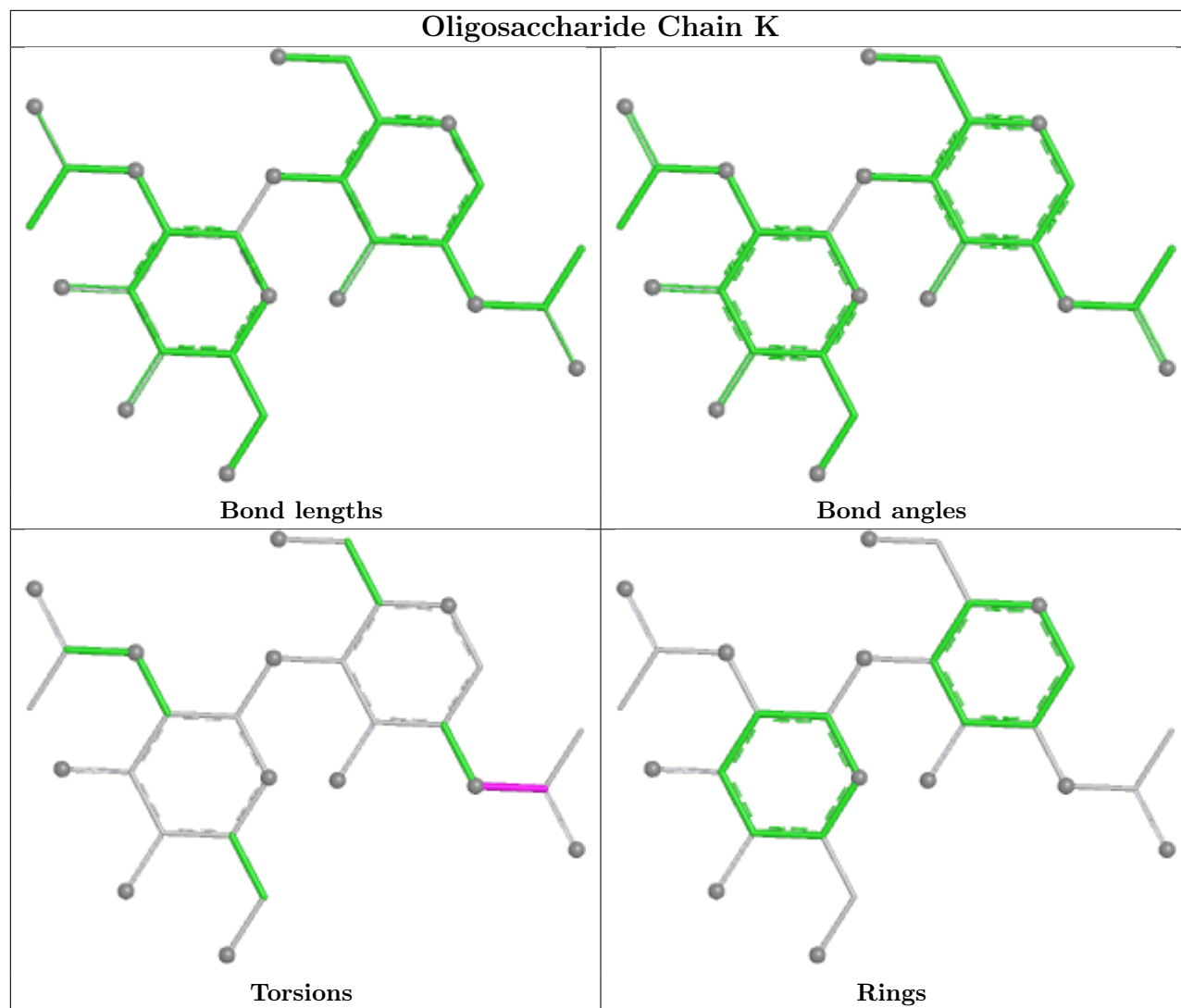


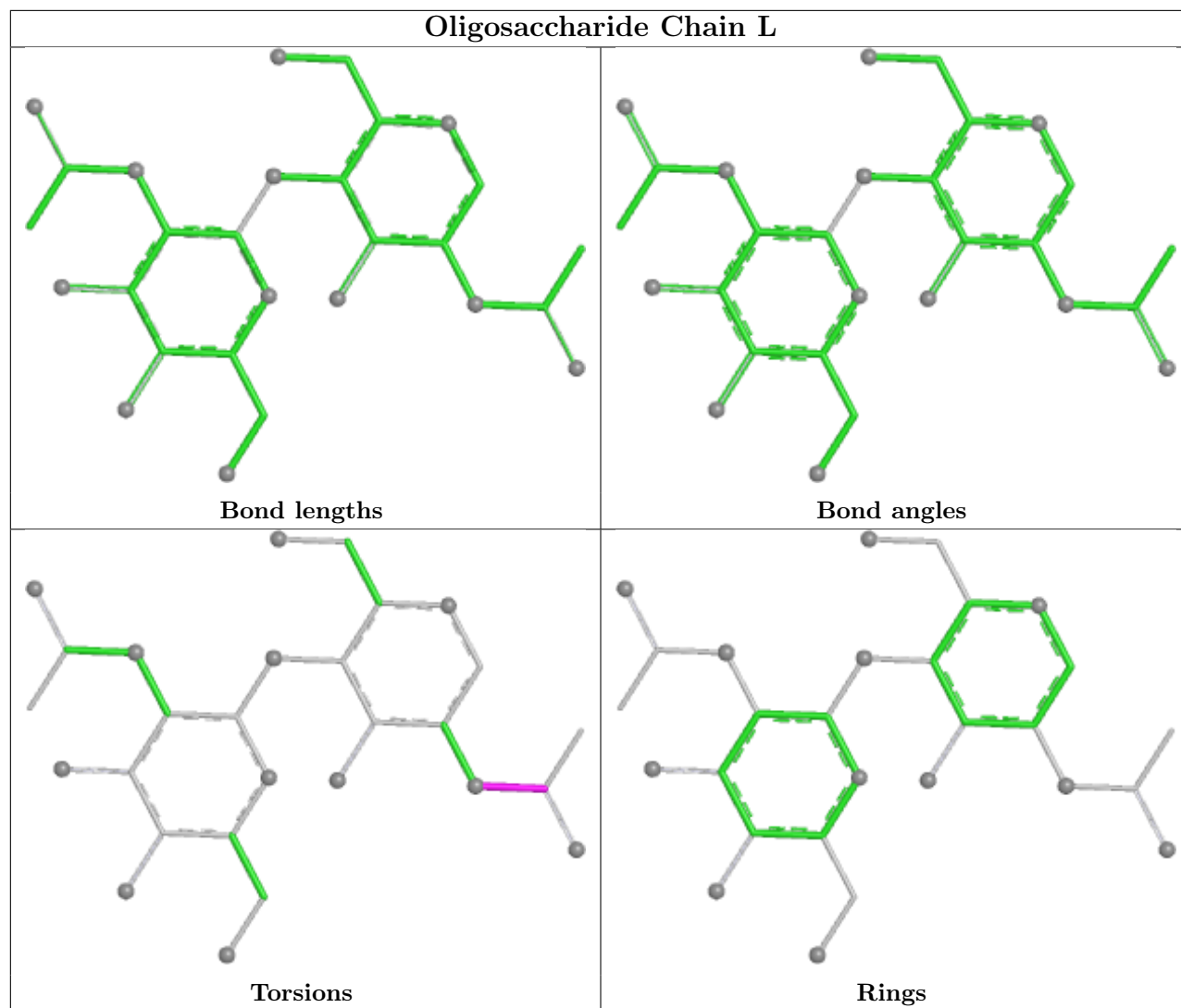


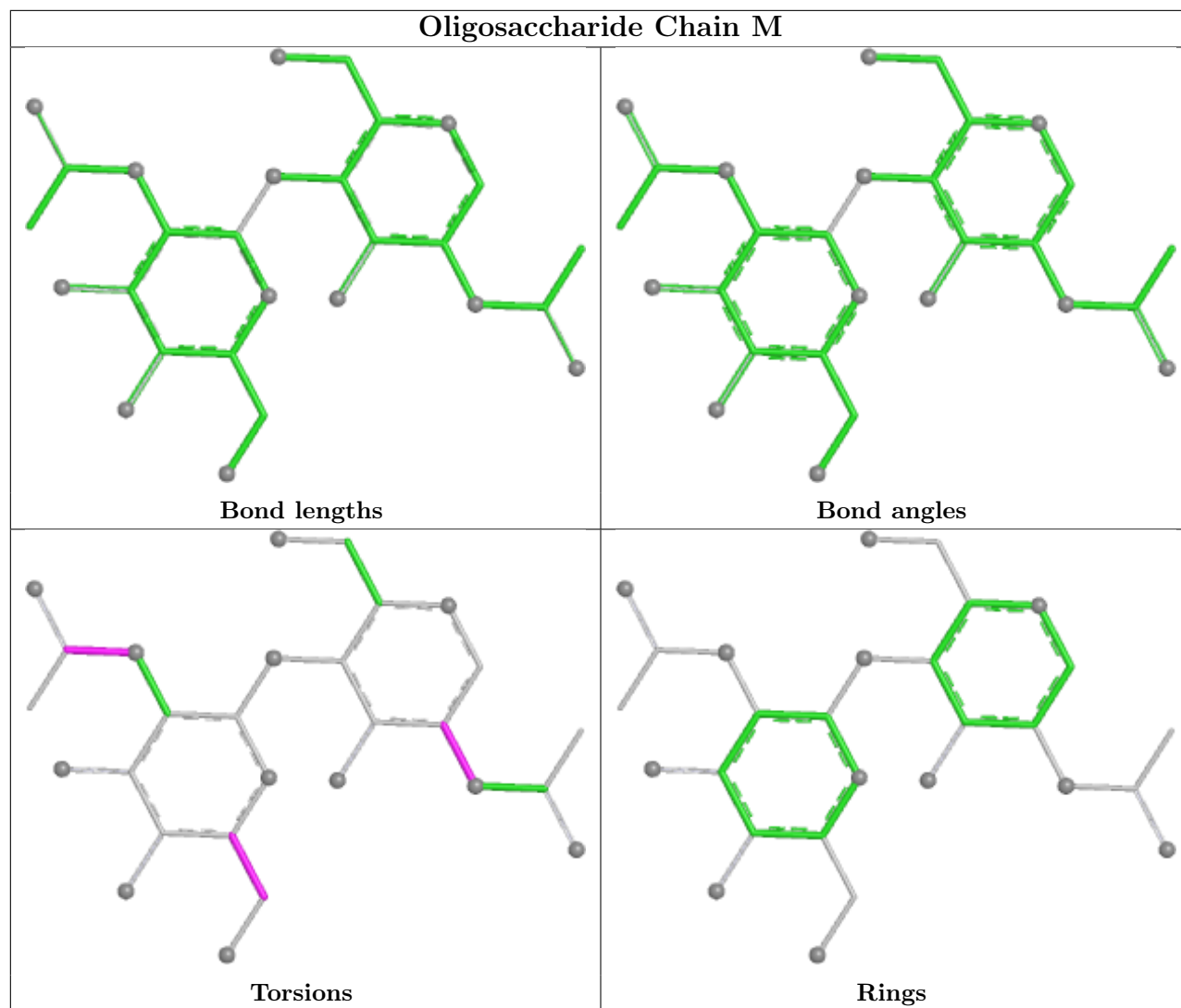


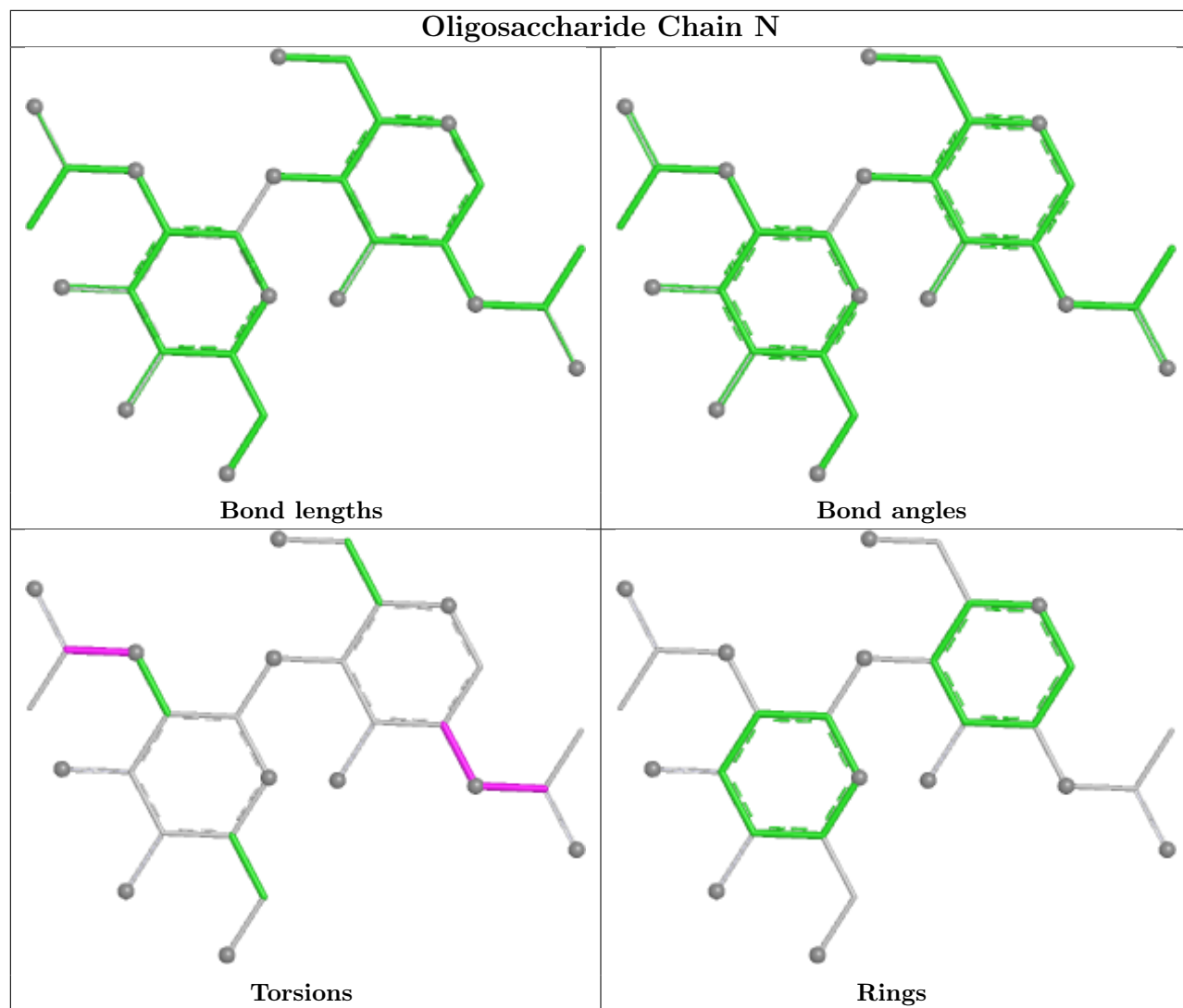


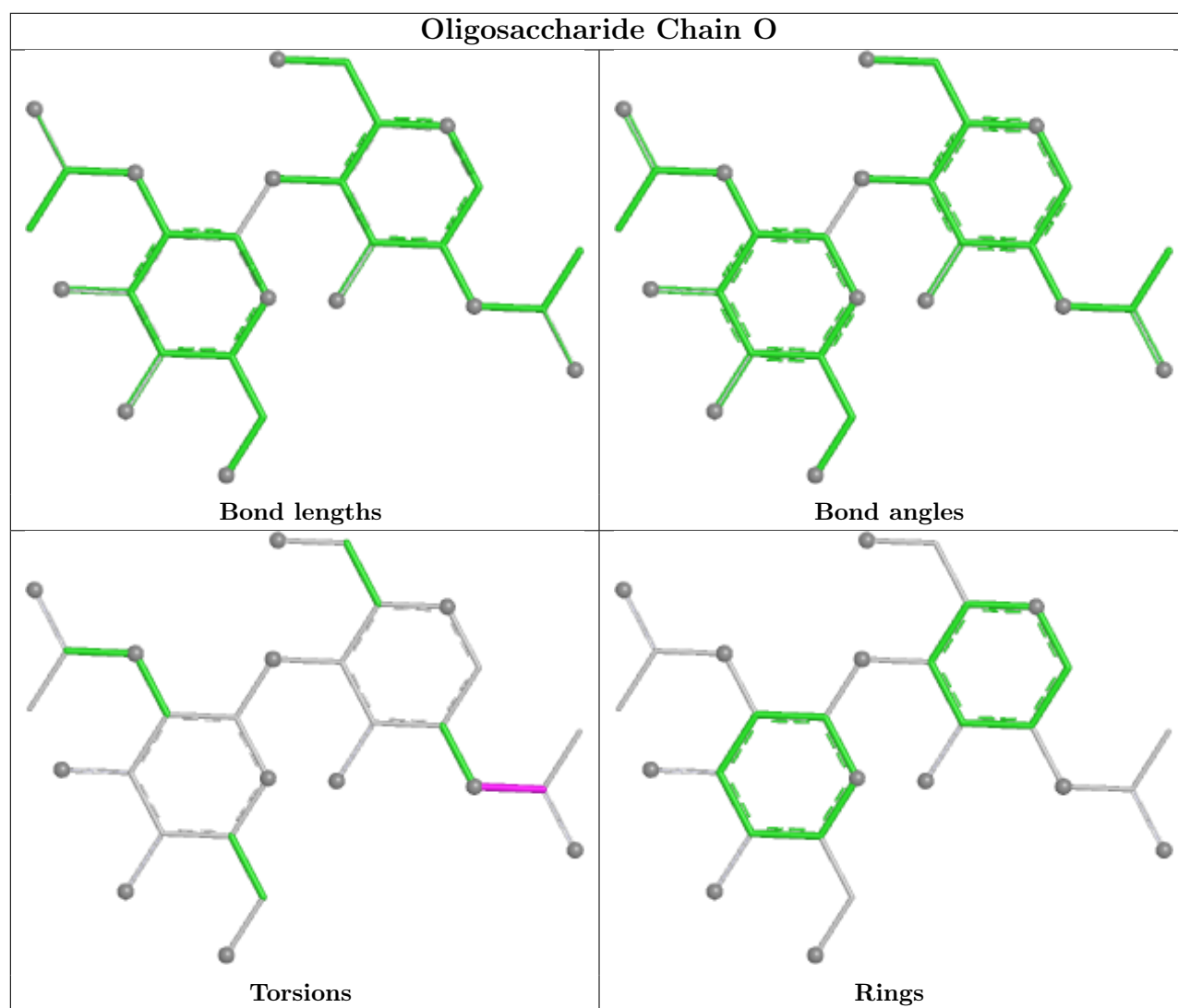


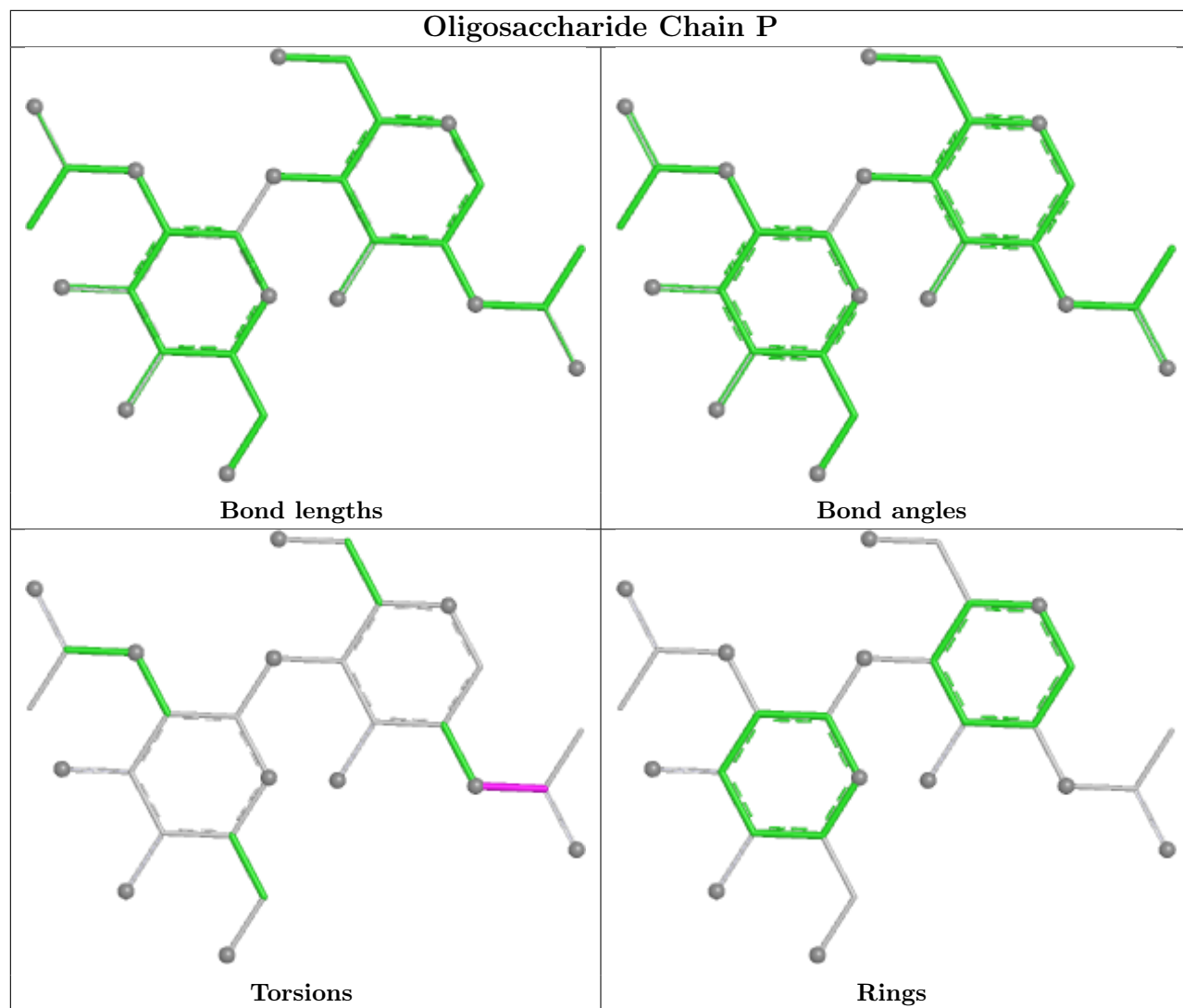


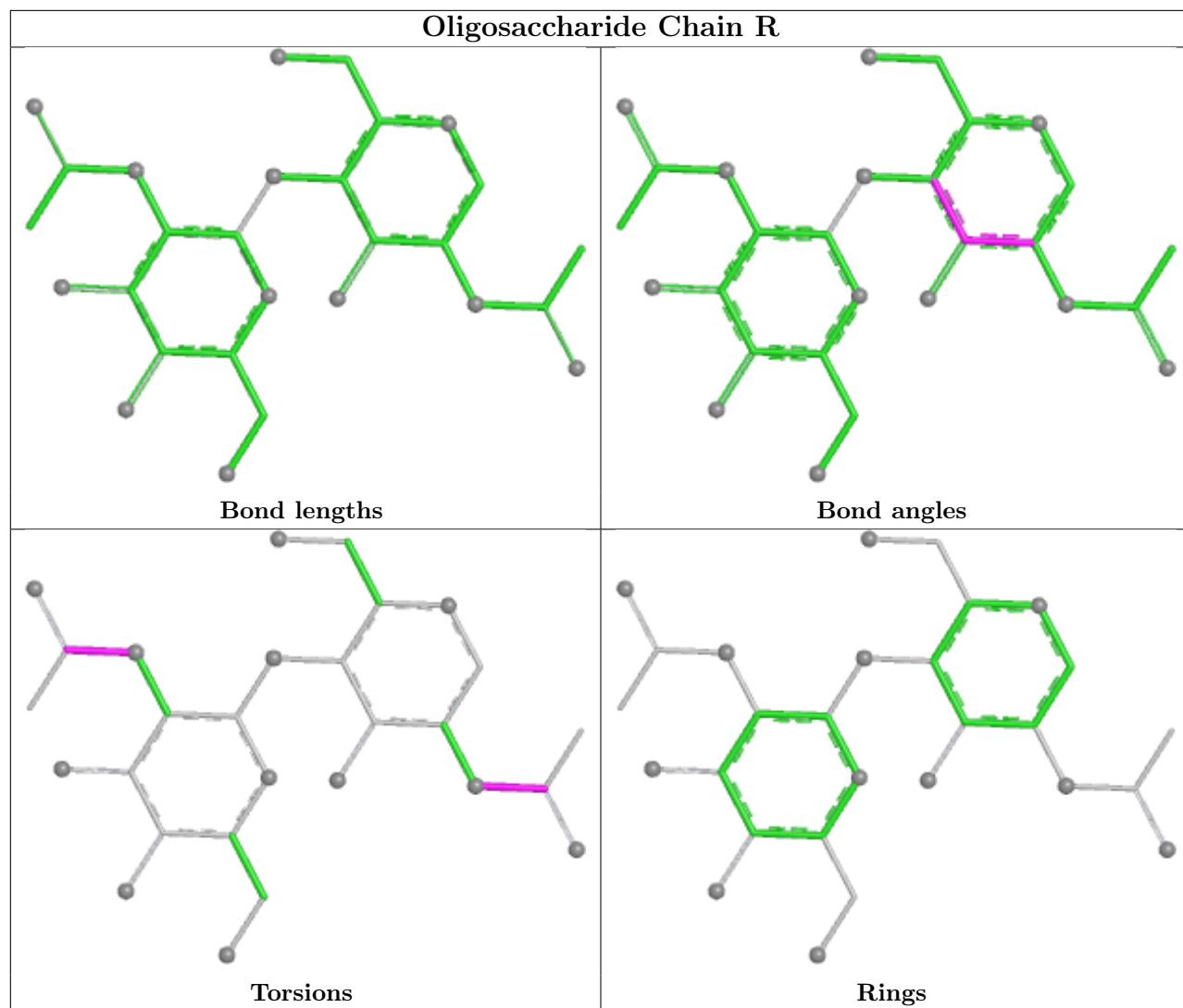


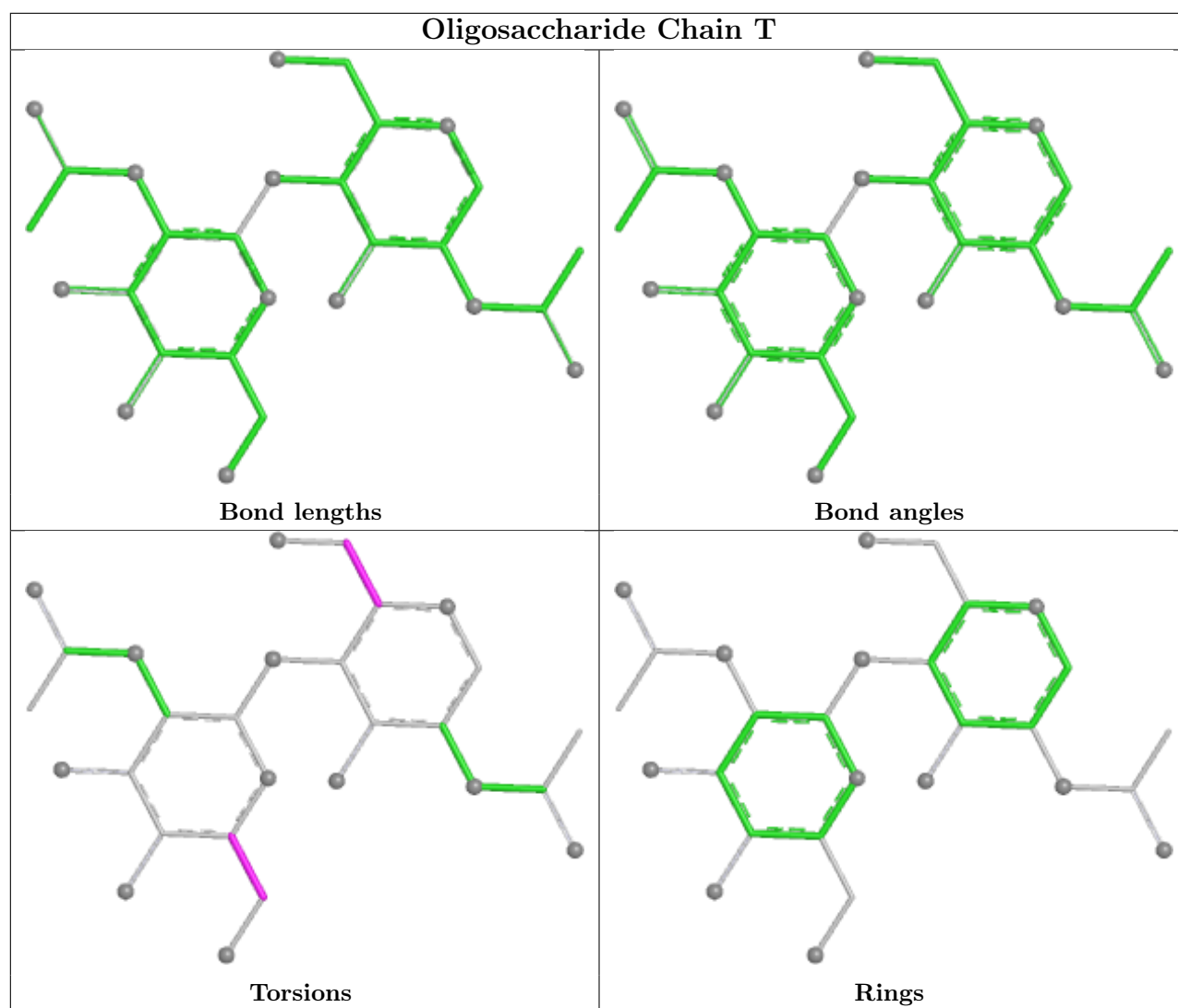


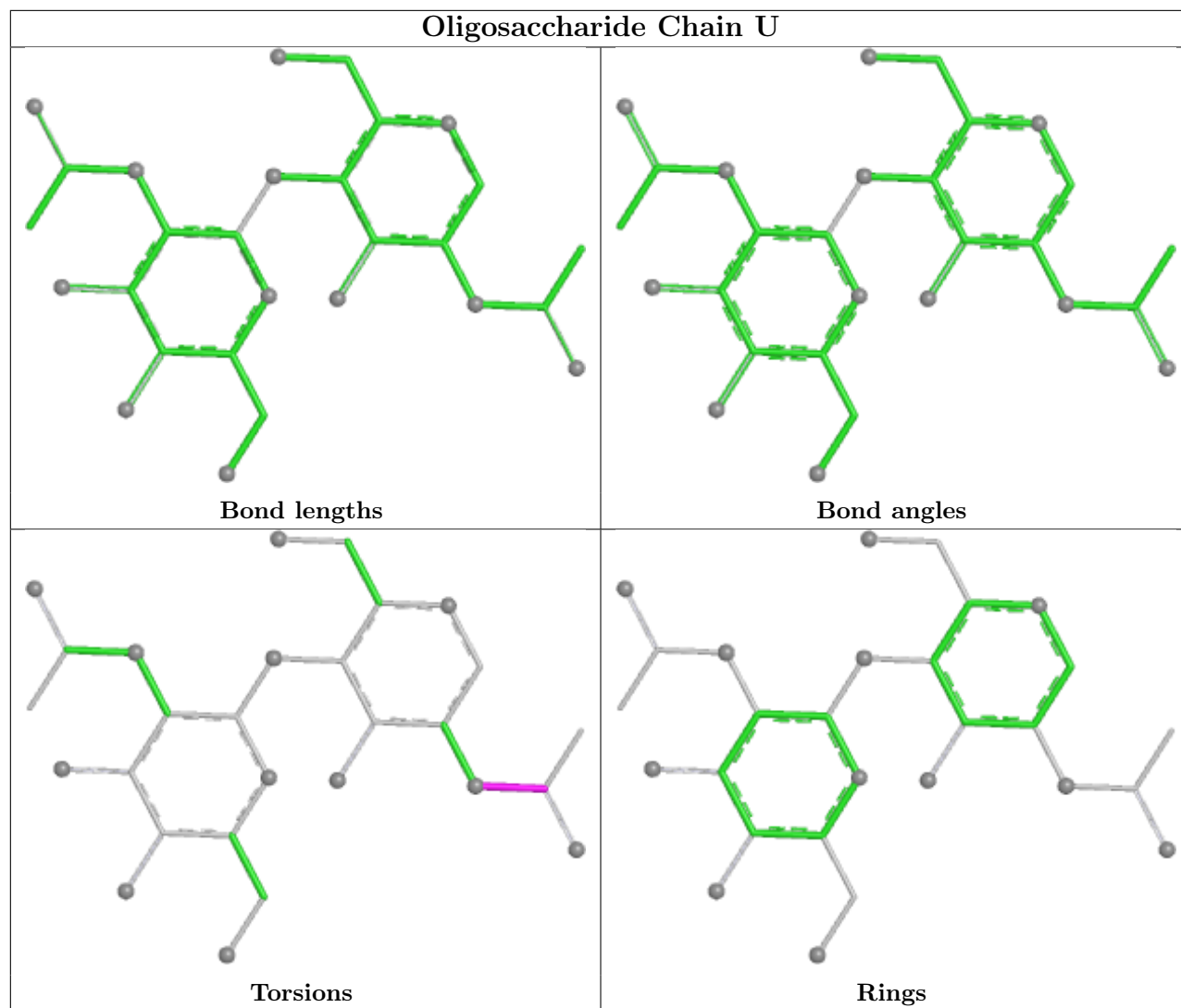


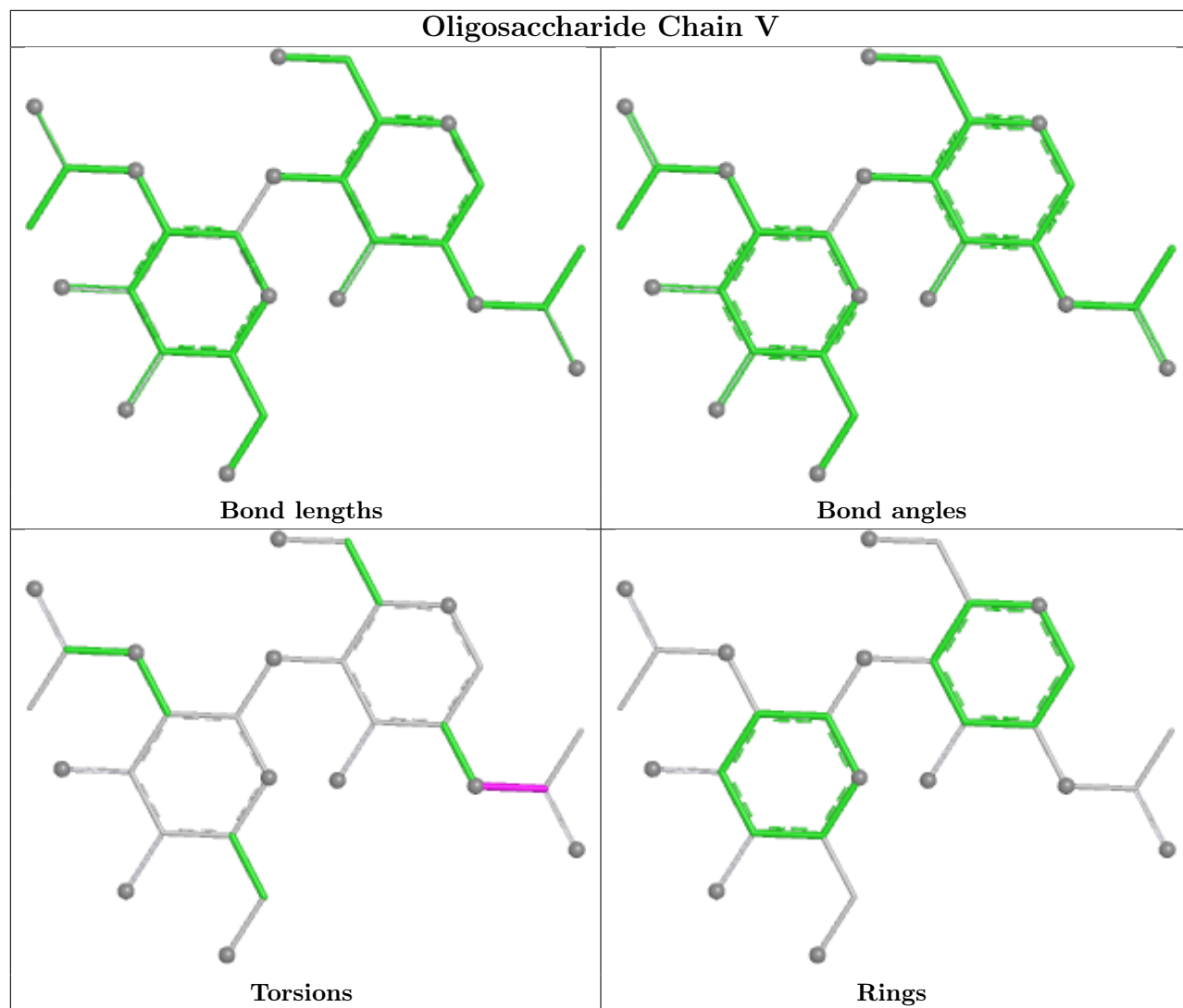


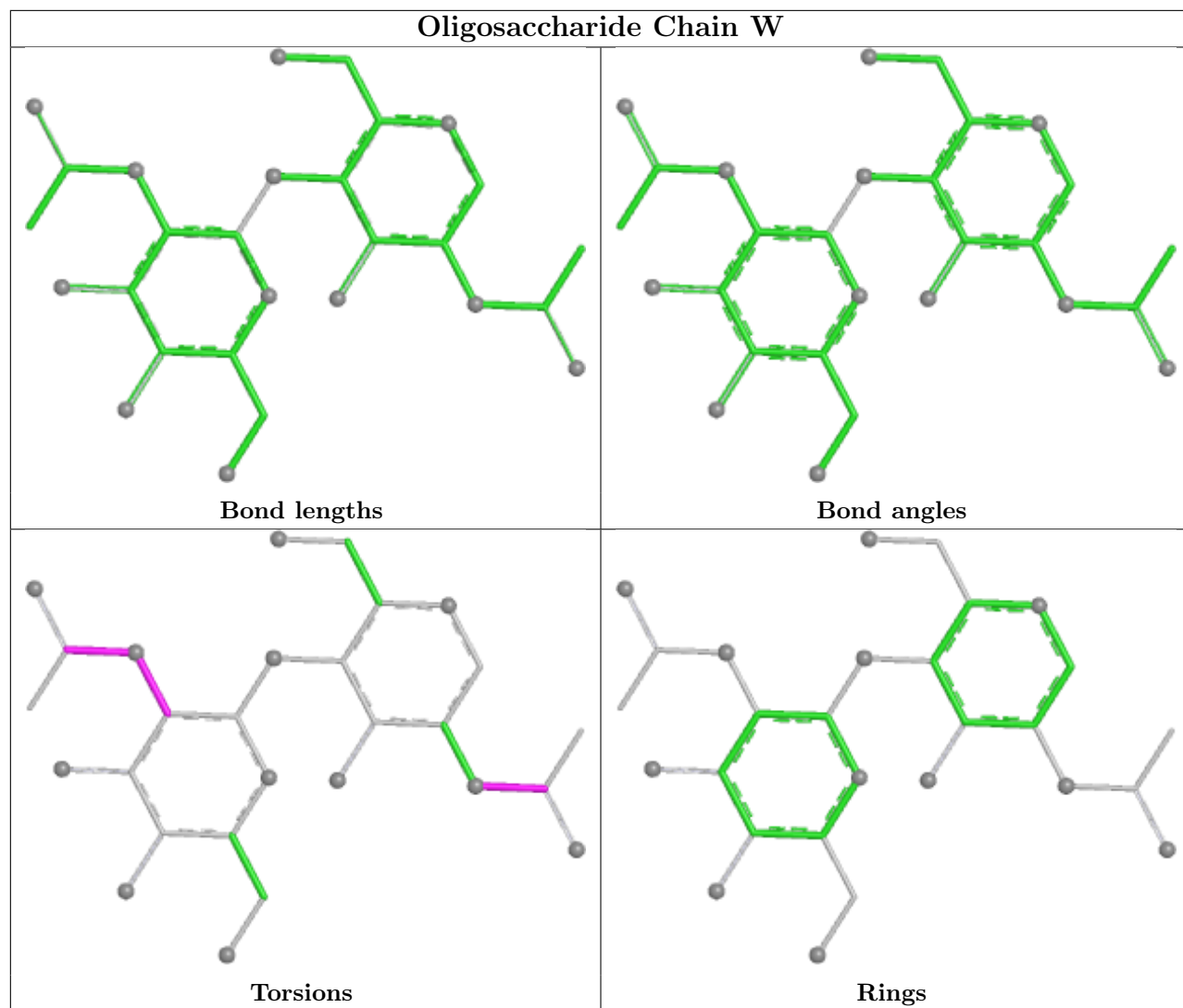


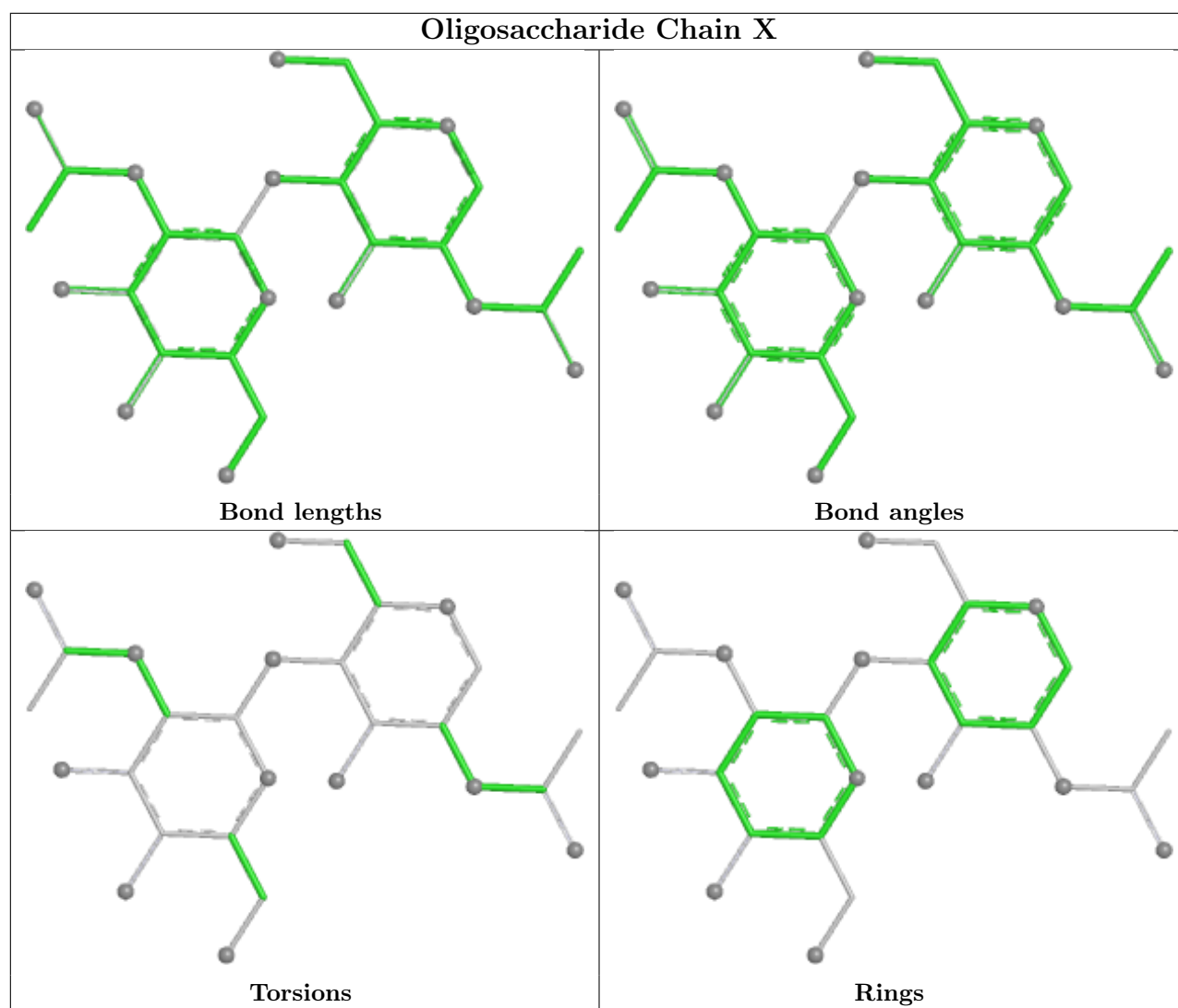


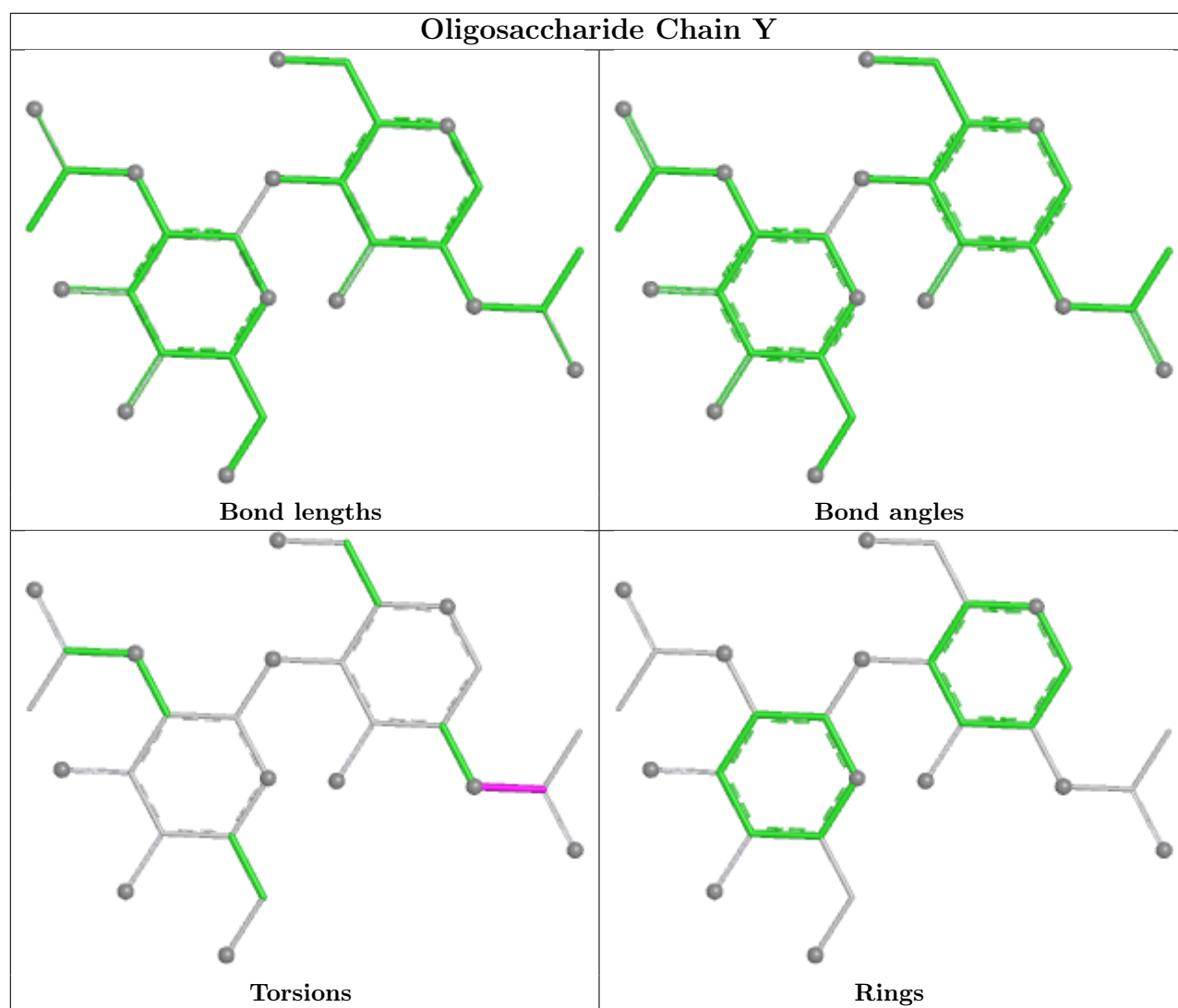


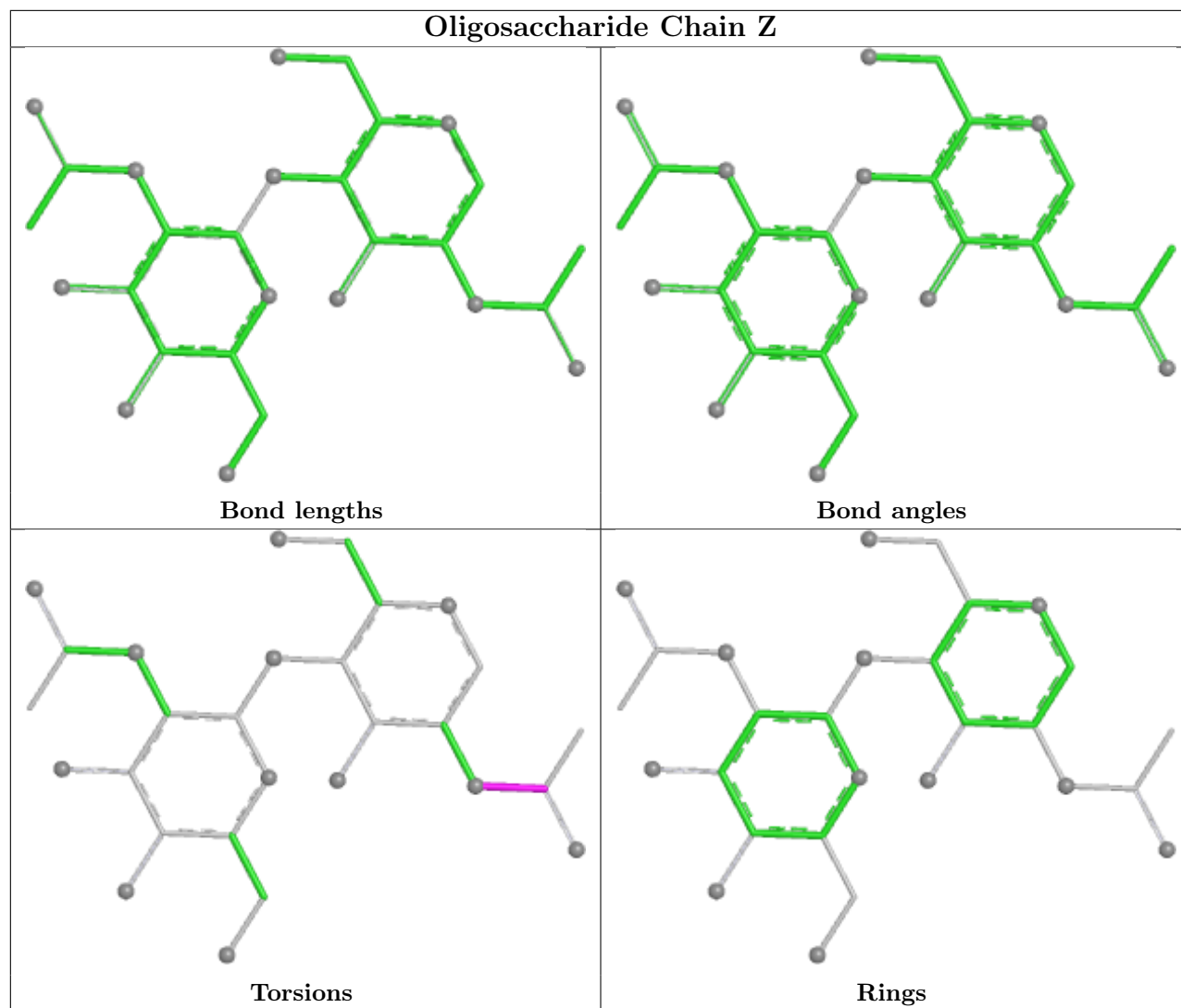


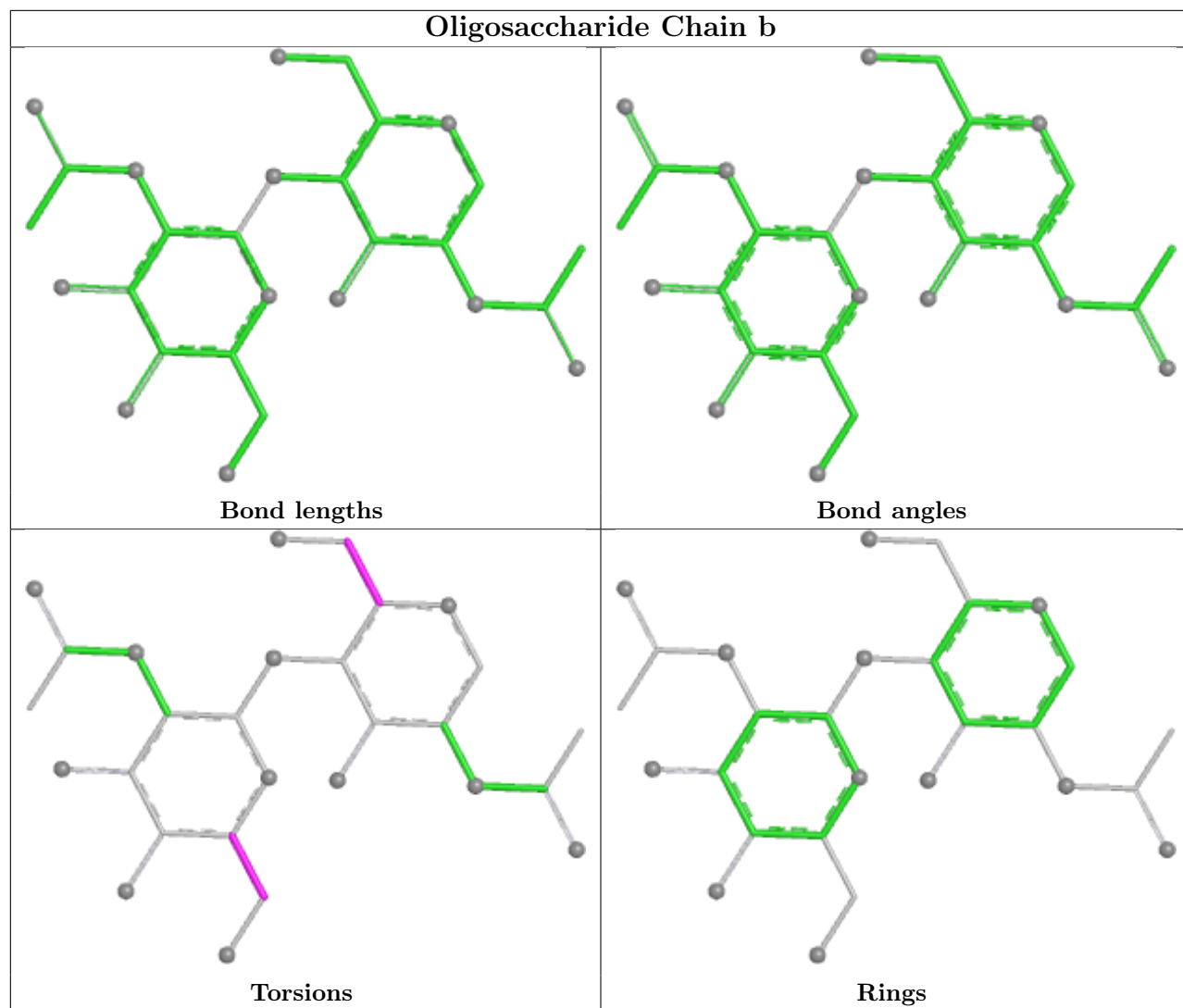


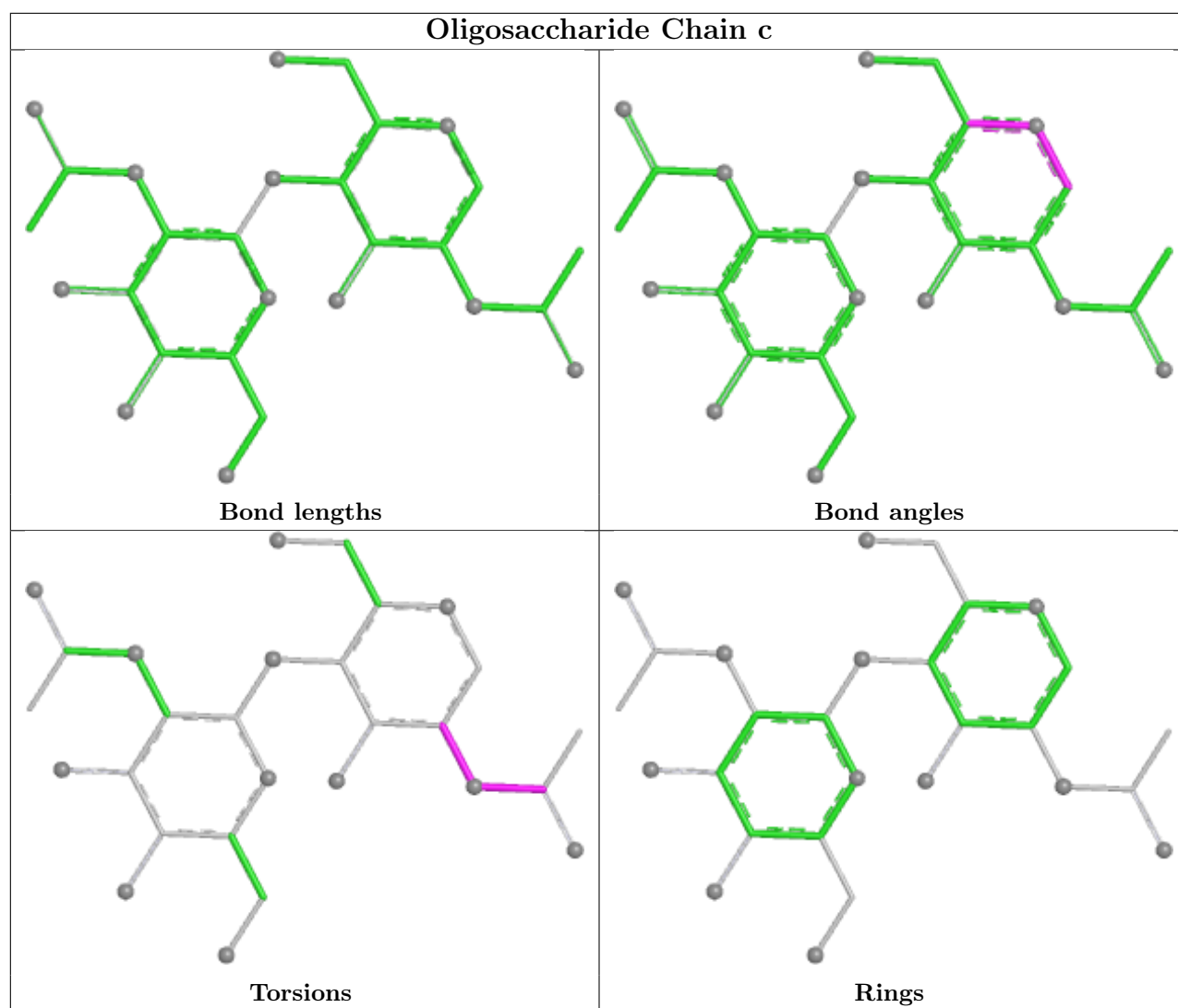


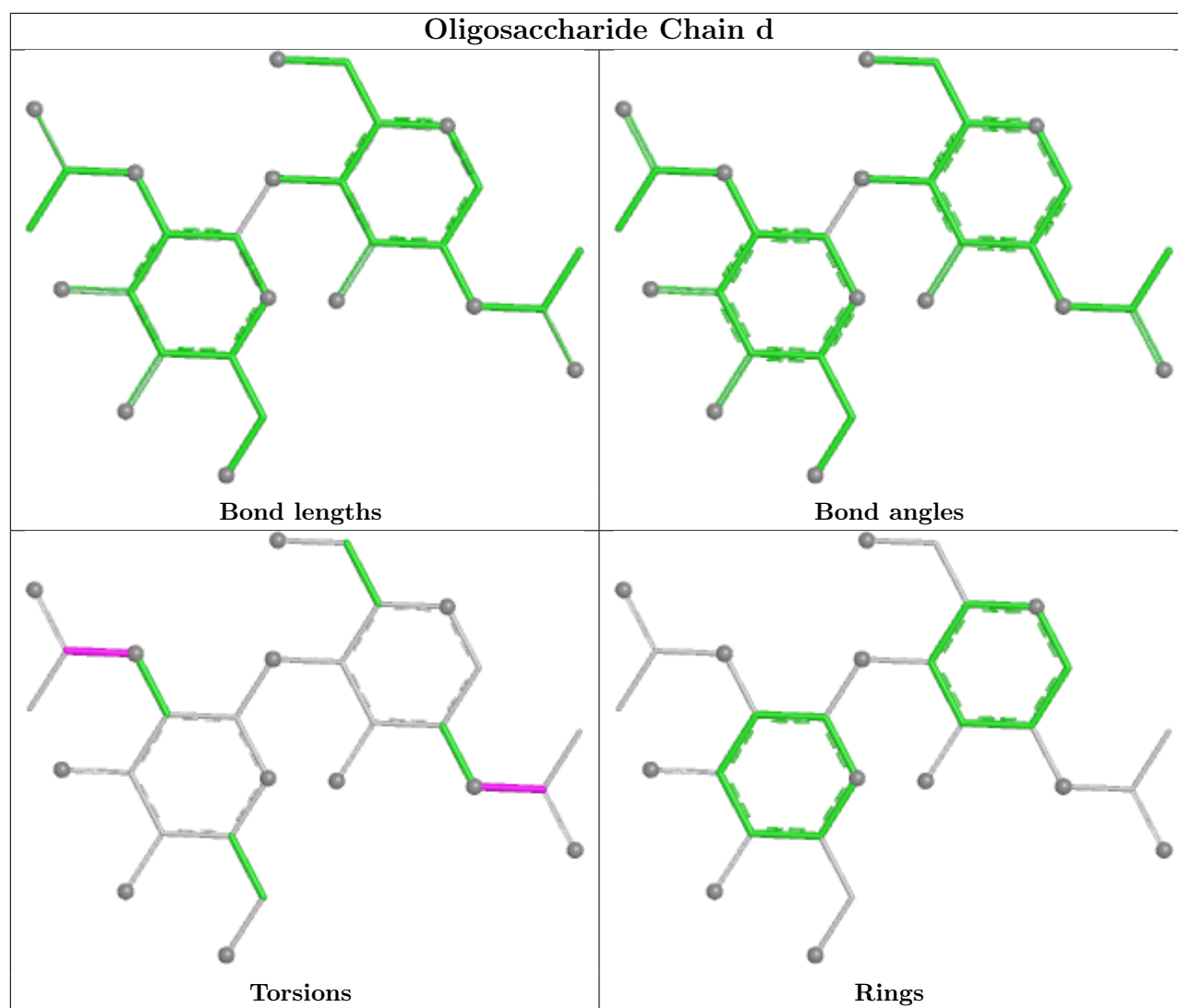


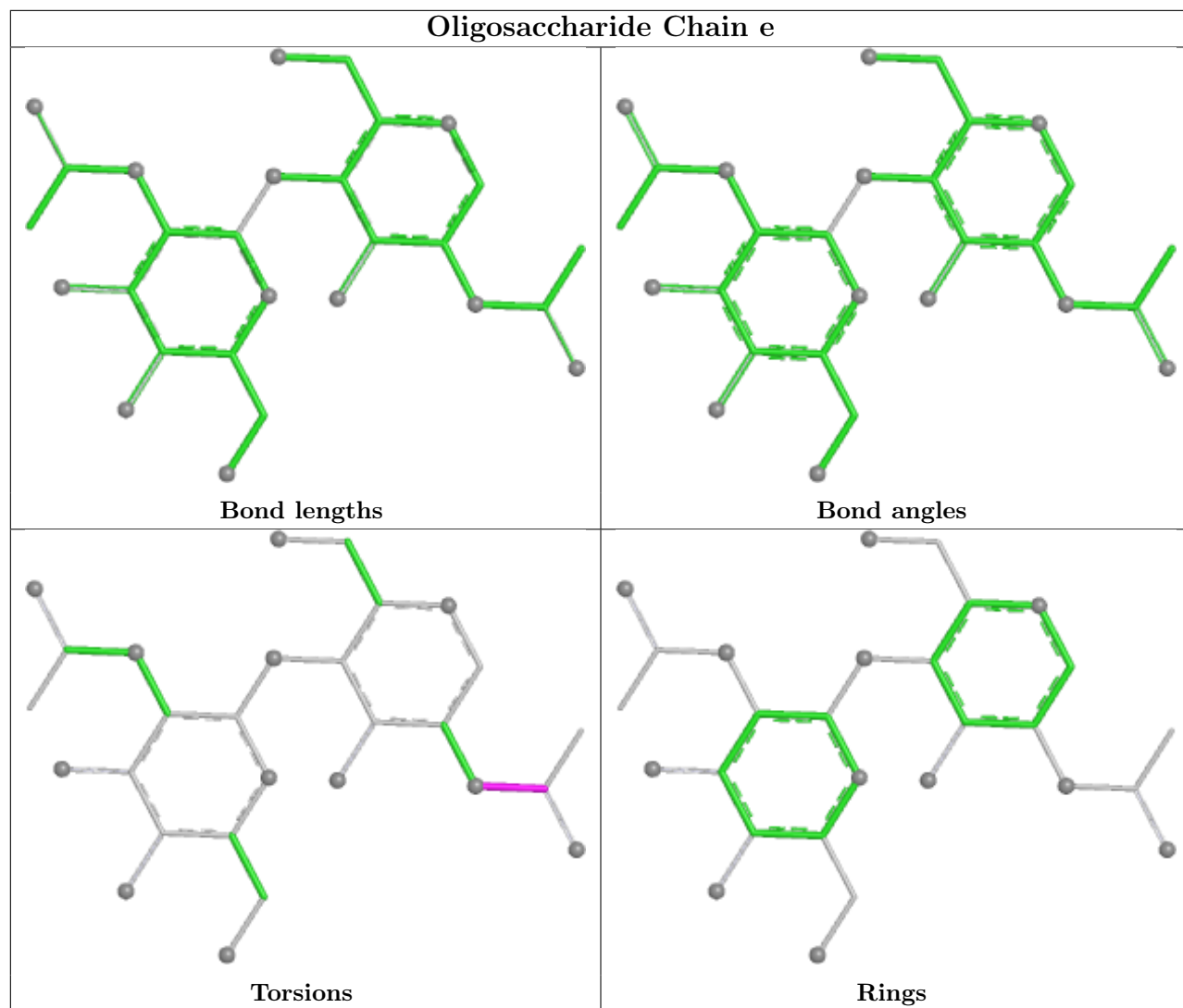


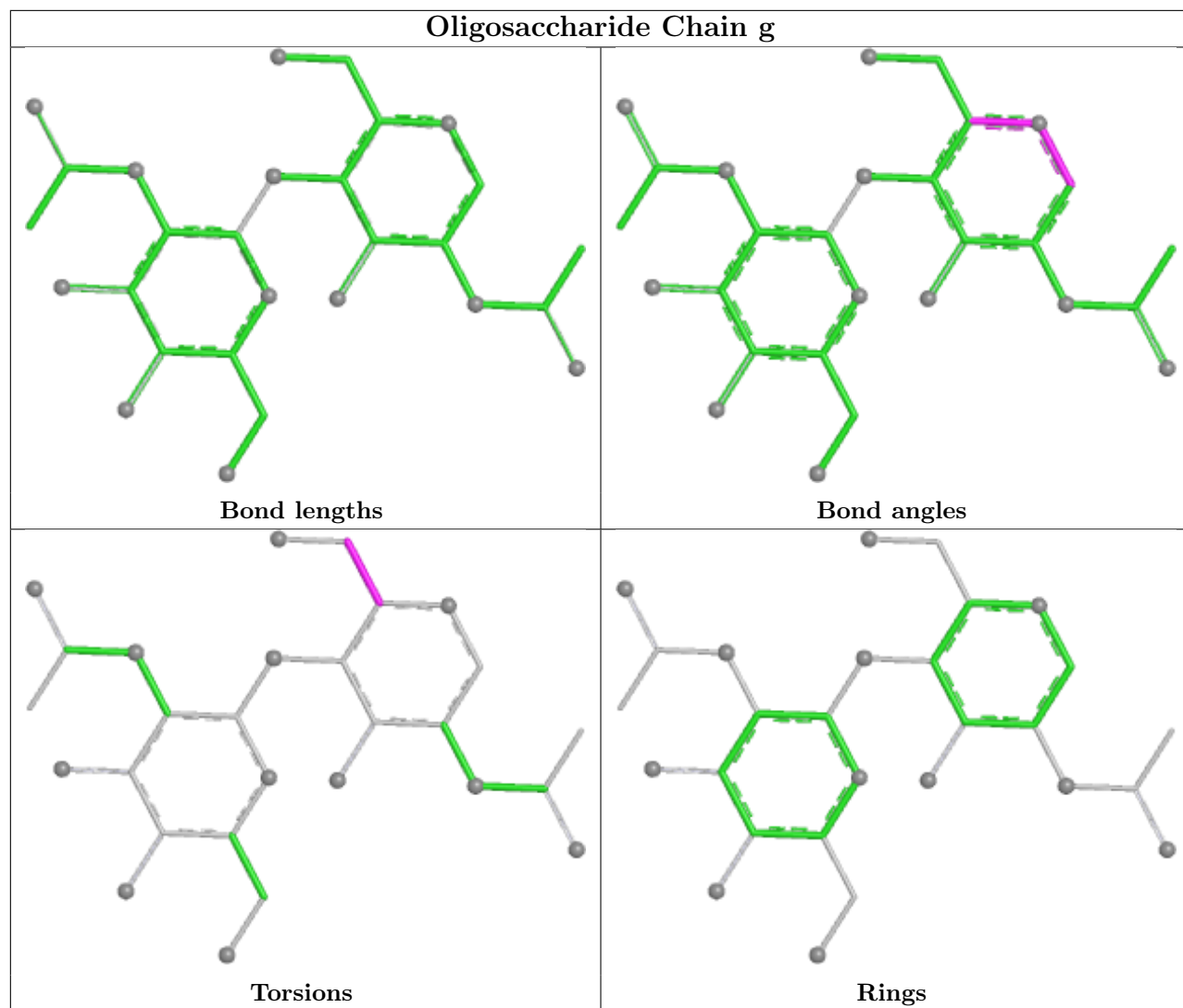


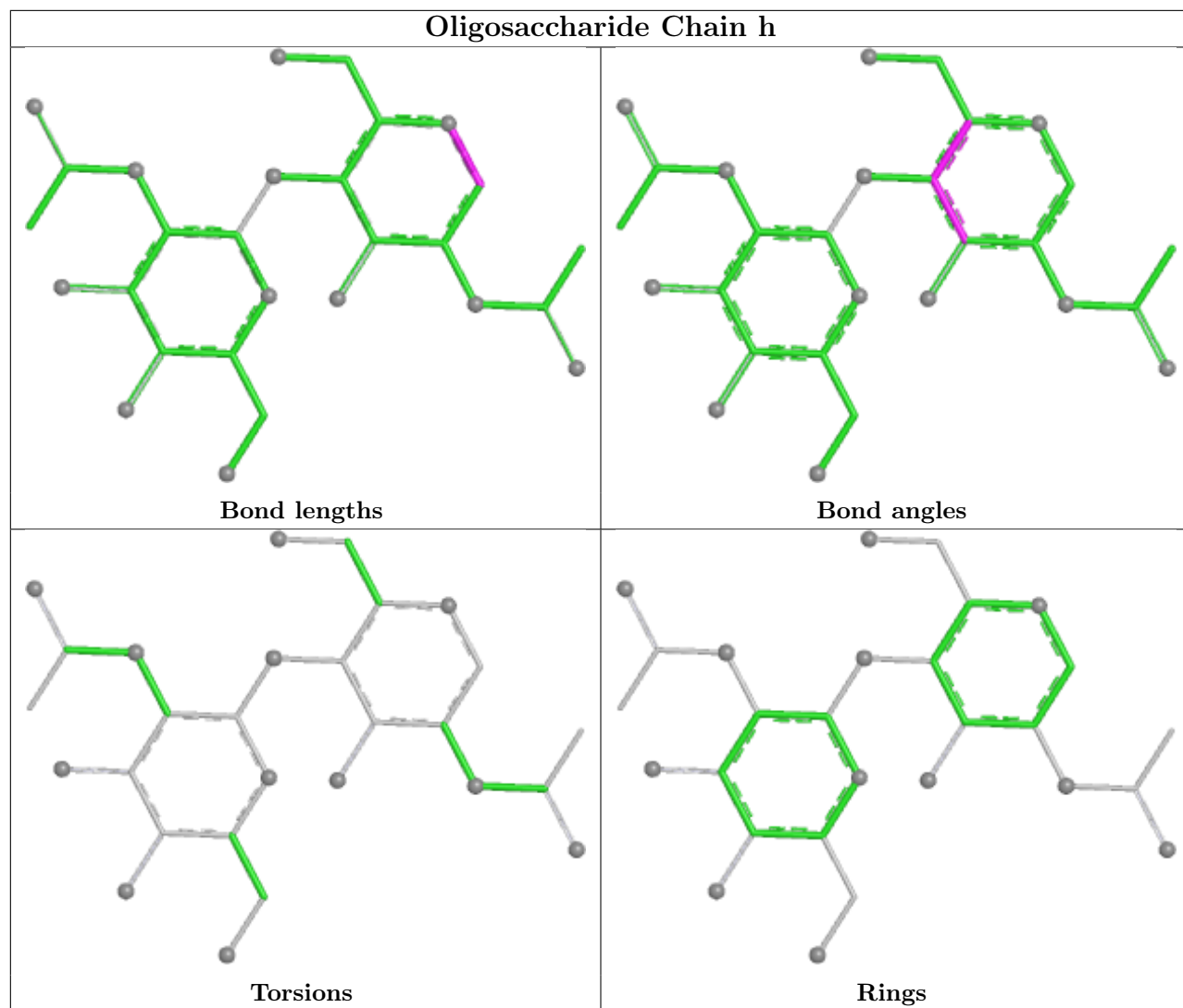


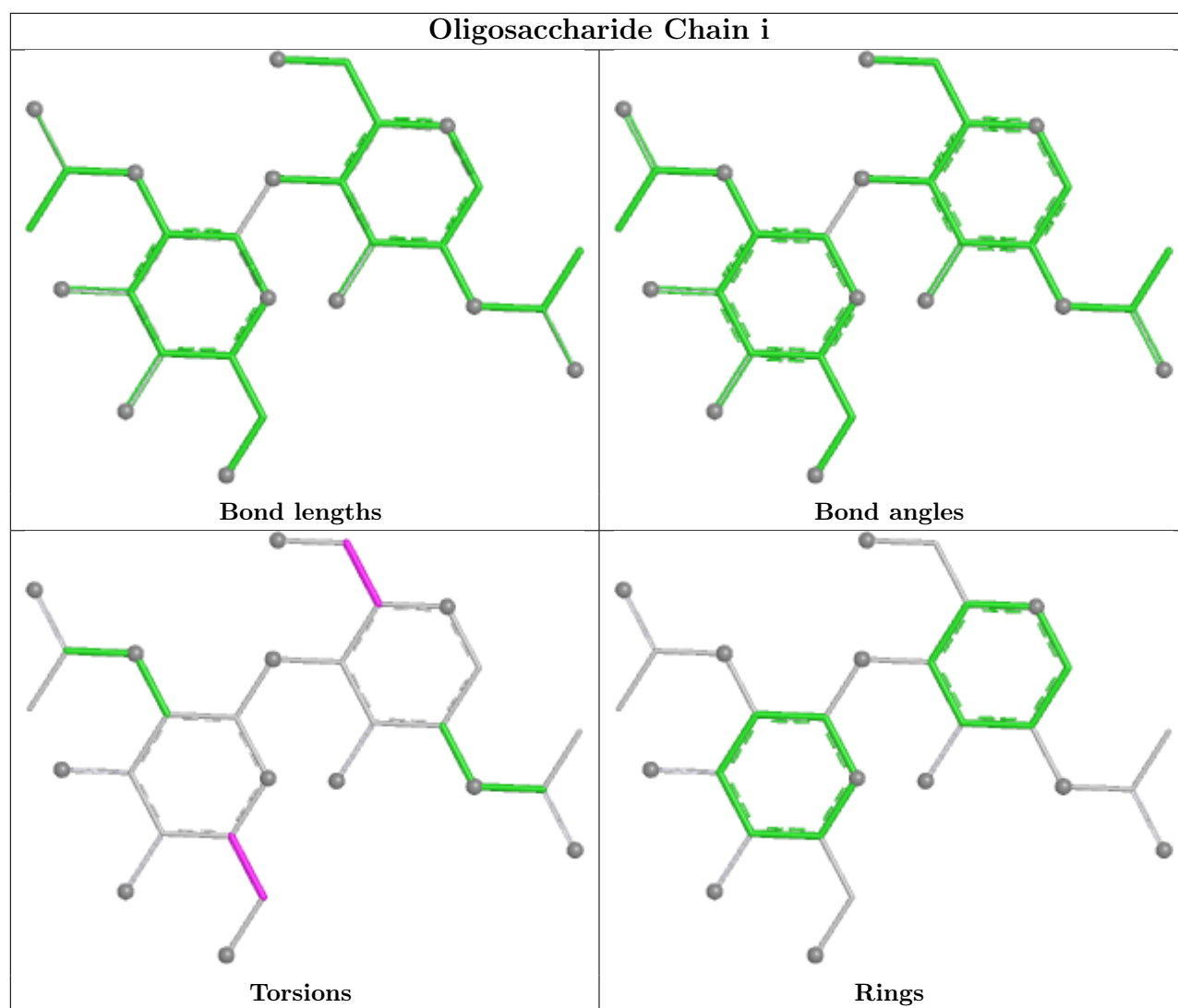


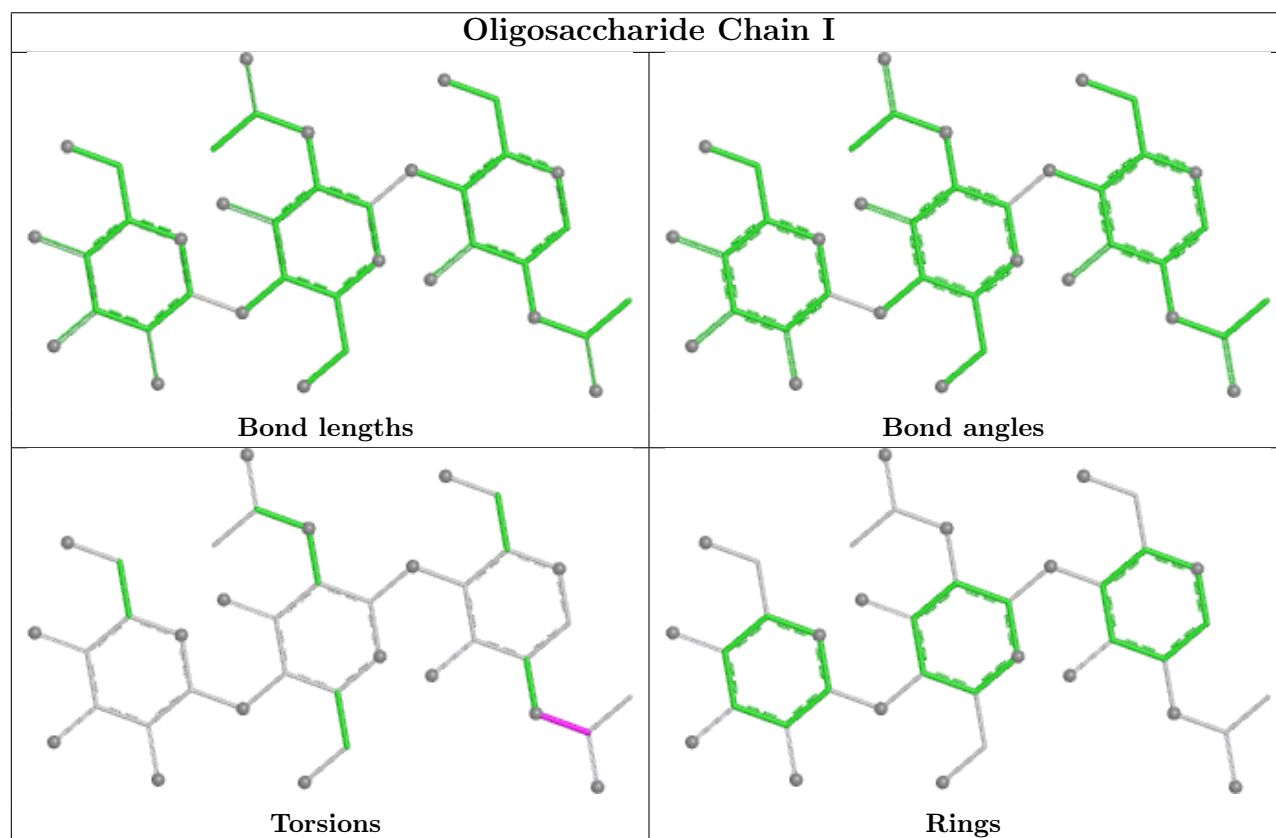
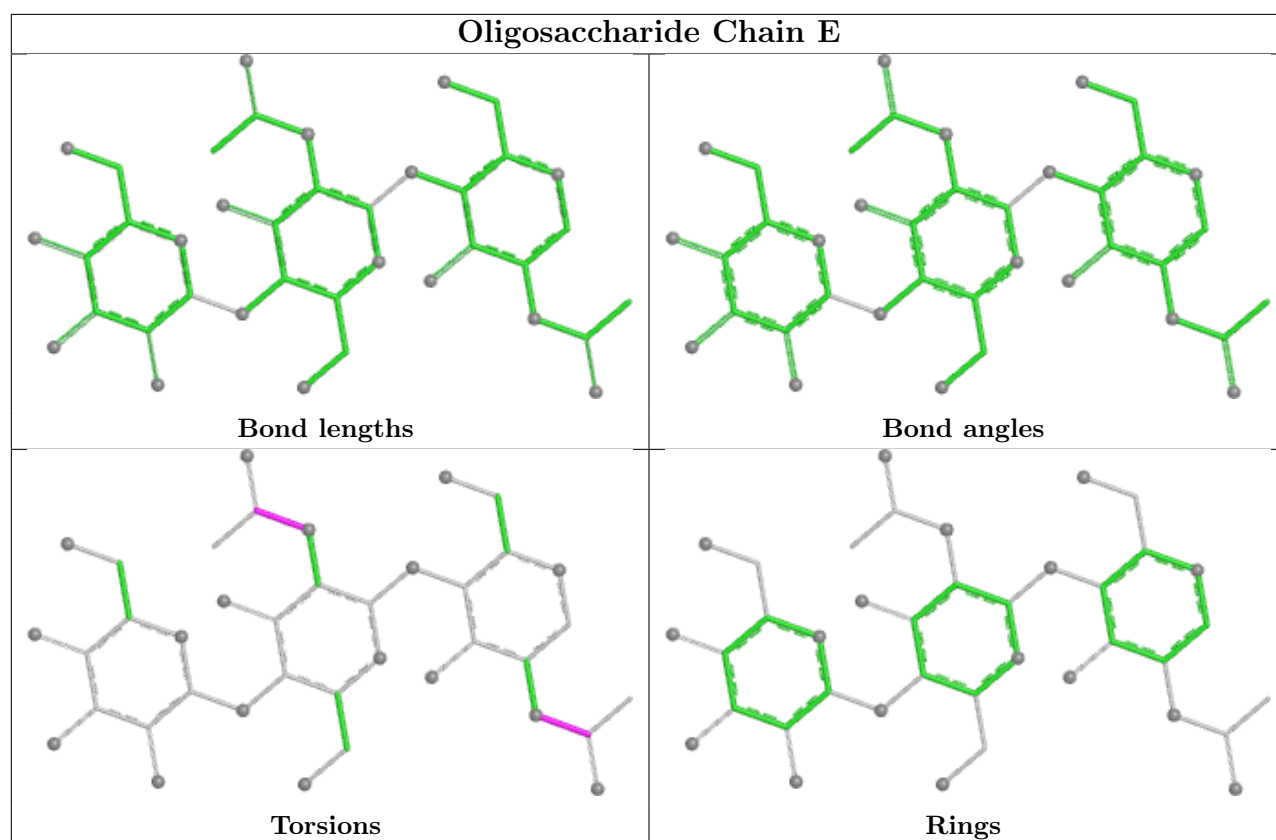


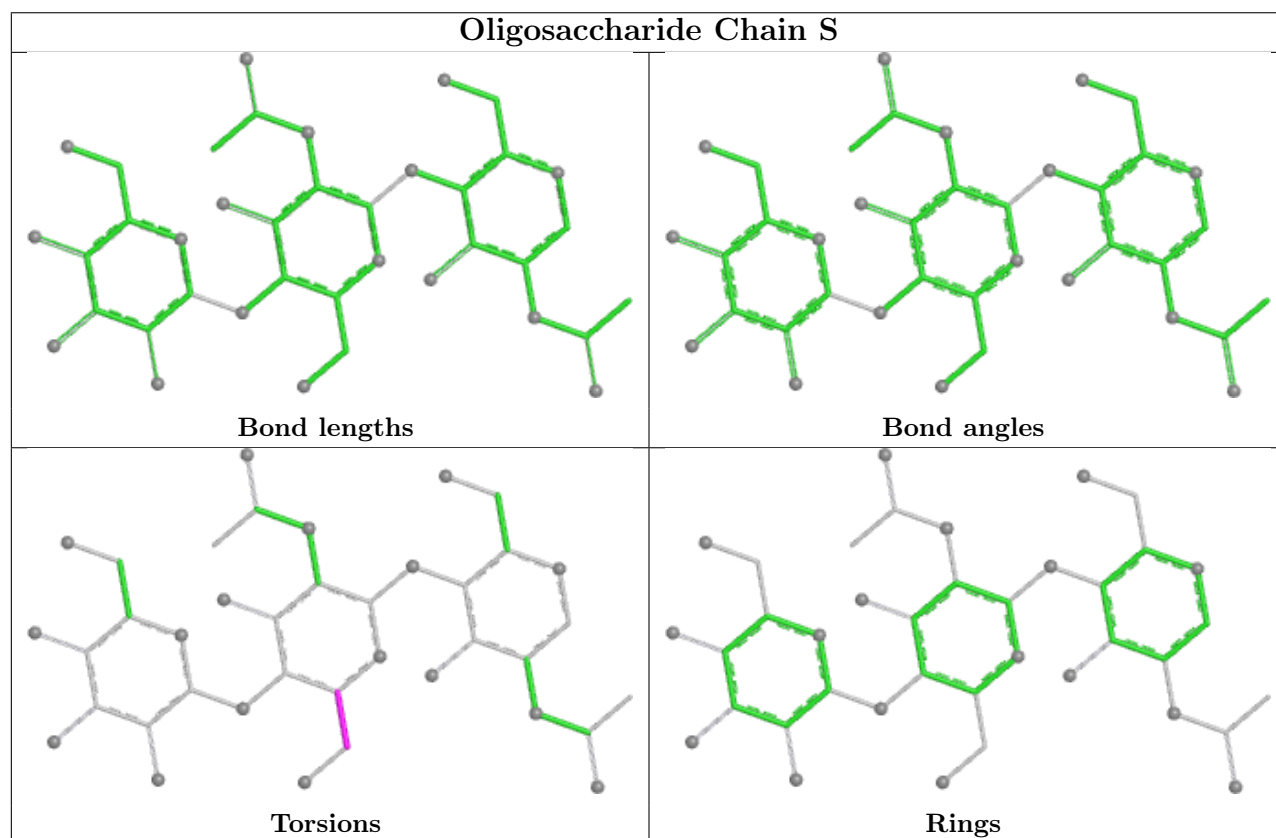
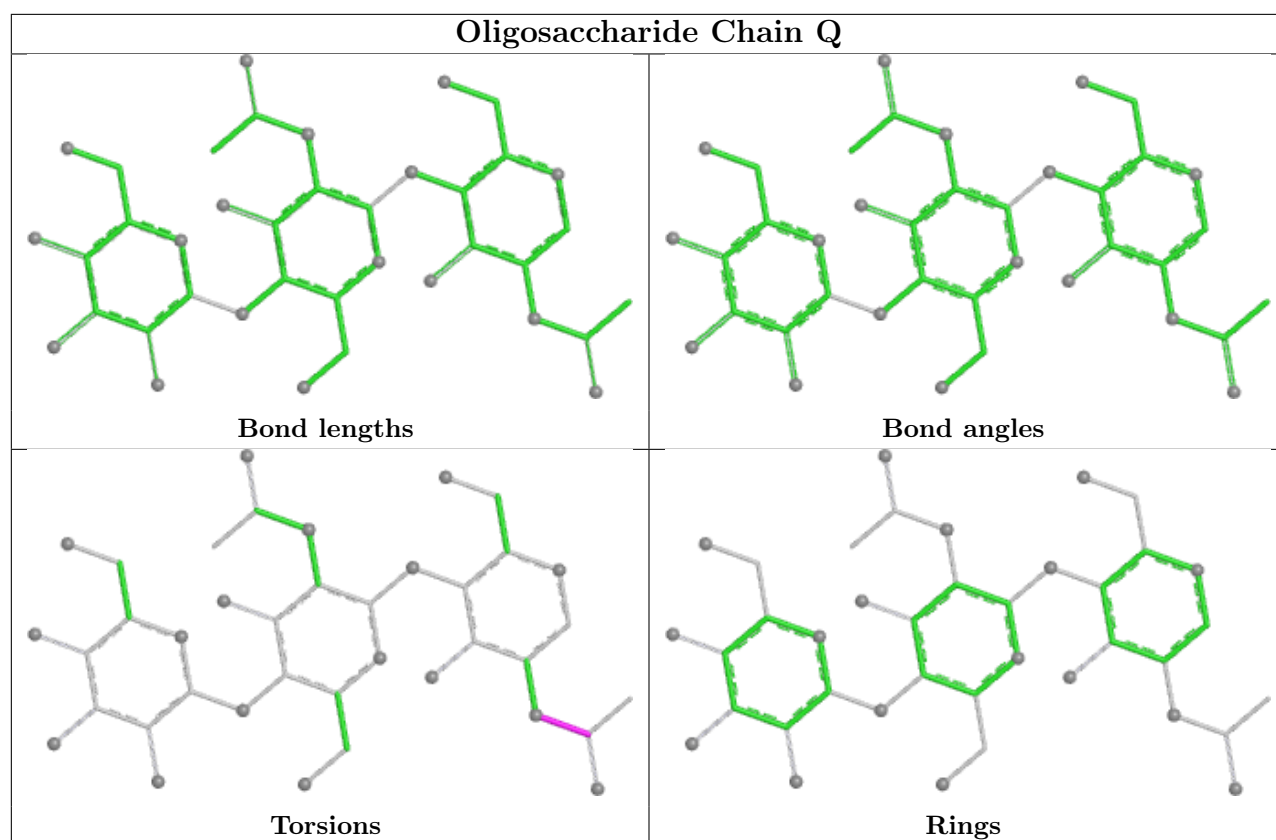


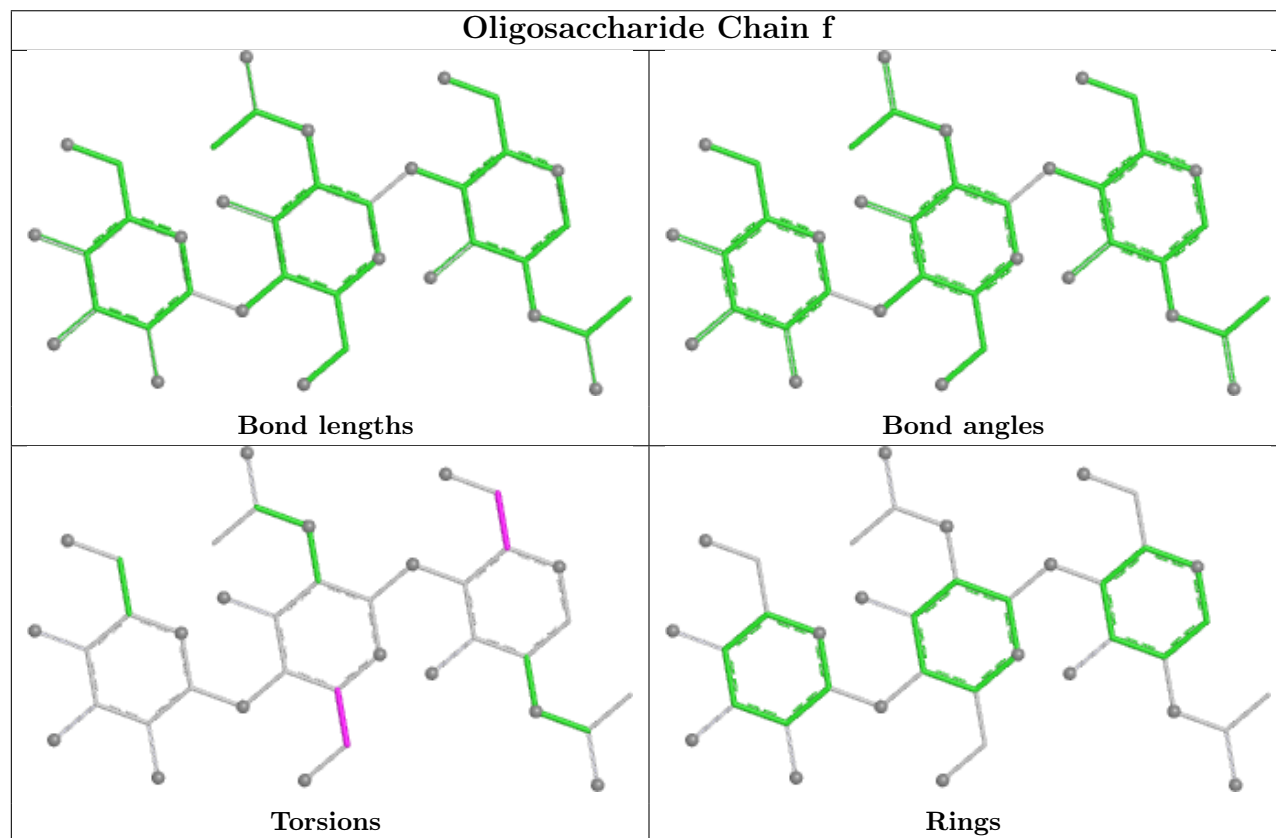
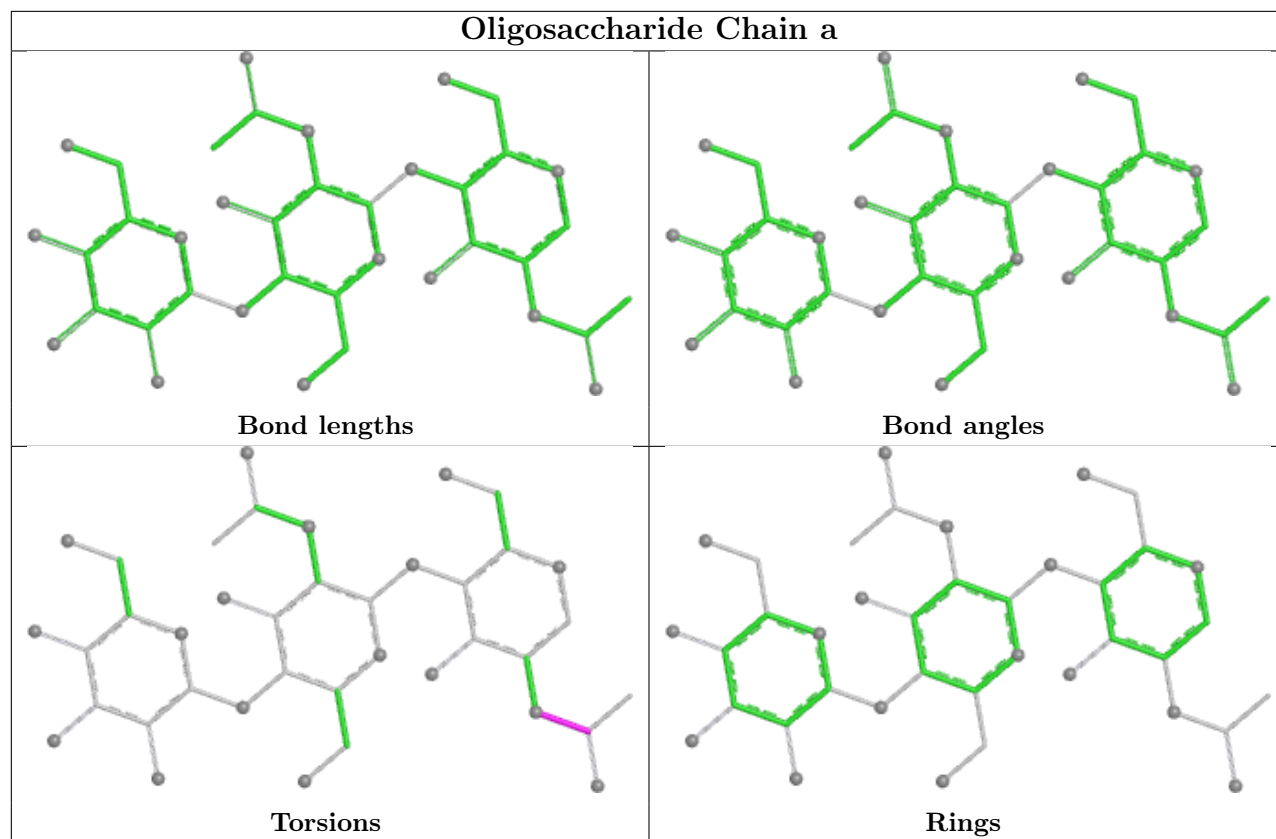


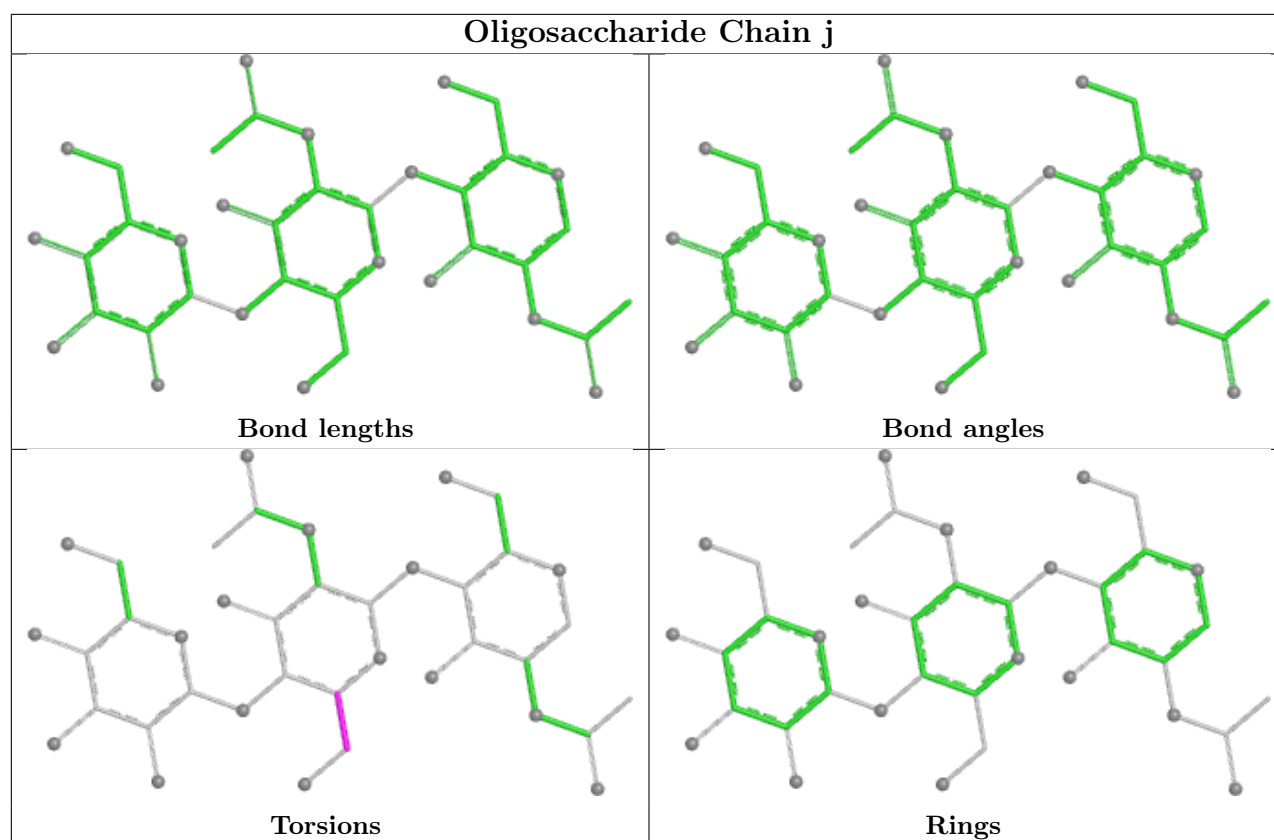












5.6 Ligand geometry [i](#)

21 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	$\# Z > 2$
4	NAG	C	1306	1	14,14,15	0.30	0	17,19,21	0.51	0
4	NAG	B	1306	1	14,14,15	0.26	0	17,19,21	0.62	0
4	NAG	B	1308	1	14,14,15	0.26	0	17,19,21	1.06	2 (11%)
4	NAG	A	1302	1	14,14,15	0.29	0	17,19,21	0.55	0
4	NAG	B	1307	1	14,14,15	0.28	0	17,19,21	0.50	0
4	NAG	C	1301	1	14,14,15	0.28	0	17,19,21	0.66	0
4	NAG	C	1302	1	14,14,15	0.28	0	17,19,21	0.57	0
4	NAG	A	1304	1	14,14,15	0.28	0	17,19,21	0.59	0
4	NAG	B	1305	1	14,14,15	0.28	0	17,19,21	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	C	1304	1	14,14,15	0.44	0	17,19,21	1.34	2 (11%)
4	NAG	B	1303	1	14,14,15	0.27	0	17,19,21	0.54	0
4	NAG	B	1304	1	14,14,15	0.45	0	17,19,21	1.35	2 (11%)
4	NAG	A	1303	1	14,14,15	0.46	0	17,19,21	1.34	2 (11%)
4	NAG	C	1303	1	14,14,15	0.29	0	17,19,21	0.58	0
4	NAG	A	1301	1	14,14,15	0.28	0	17,19,21	0.62	0
4	NAG	A	1305	1	14,14,15	0.27	0	17,19,21	0.58	0
4	NAG	B	1301	1	14,14,15	0.28	0	17,19,21	0.63	0
4	NAG	B	1302	1	14,14,15	0.28	0	17,19,21	0.57	0
4	NAG	C	1305	1	14,14,15	0.28	0	17,19,21	0.59	0
4	NAG	A	1306	1	14,14,15	0.28	0	17,19,21	0.59	0
4	NAG	A	1307	1	14,14,15	0.28	0	17,19,21	0.90	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	C	1306	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1306	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1308	1	-	4/6/23/26	0/1/1/1
4	NAG	A	1302	1	-	3/6/23/26	0/1/1/1
4	NAG	B	1307	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1301	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1302	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1304	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1305	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1304	1	-	4/6/23/26	0/1/1/1
4	NAG	B	1303	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1304	1	-	4/6/23/26	0/1/1/1
4	NAG	A	1303	1	-	4/6/23/26	0/1/1/1
4	NAG	C	1303	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1301	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1305	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1301	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1302	1	-	3/6/23/26	0/1/1/1
4	NAG	C	1305	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1306	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1307	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1303	NAG	C2-N2-C7	4.61	129.07	122.90
4	B	1304	NAG	C2-N2-C7	4.59	129.05	122.90
4	C	1304	NAG	C2-N2-C7	4.55	129.00	122.90
4	B	1308	NAG	C1-O5-C5	2.79	115.92	112.19
4	B	1308	NAG	C2-N2-C7	2.45	126.18	122.90
4	C	1304	NAG	C1-C2-N2	2.16	113.84	110.43
4	B	1304	NAG	C1-C2-N2	2.15	113.82	110.43
4	A	1303	NAG	C1-C2-N2	2.10	113.74	110.43

There are no chirality outliers.

All (50) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1306	NAG	C8-C7-N2-C2
4	A	1306	NAG	O7-C7-N2-C2
4	B	1305	NAG	C8-C7-N2-C2
4	B	1305	NAG	O7-C7-N2-C2
4	B	1308	NAG	C1-C2-N2-C7
4	B	1308	NAG	C8-C7-N2-C2
4	B	1308	NAG	O7-C7-N2-C2
4	A	1304	NAG	O7-C7-N2-C2
4	A	1307	NAG	O7-C7-N2-C2
4	A	1304	NAG	C8-C7-N2-C2
4	A	1307	NAG	C8-C7-N2-C2
4	B	1301	NAG	C8-C7-N2-C2
4	B	1301	NAG	O7-C7-N2-C2
4	B	1302	NAG	C8-C7-N2-C2
4	C	1301	NAG	C8-C7-N2-C2
4	C	1301	NAG	O7-C7-N2-C2
4	C	1302	NAG	C8-C7-N2-C2
4	A	1303	NAG	C8-C7-N2-C2
4	A	1303	NAG	O7-C7-N2-C2
4	B	1302	NAG	O7-C7-N2-C2
4	B	1304	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
4	B	1304	NAG	O7-C7-N2-C2
4	C	1302	NAG	O7-C7-N2-C2
4	C	1304	NAG	C8-C7-N2-C2
4	C	1304	NAG	O7-C7-N2-C2
4	C	1305	NAG	C8-C7-N2-C2
4	C	1305	NAG	O7-C7-N2-C2
4	A	1301	NAG	C8-C7-N2-C2
4	A	1302	NAG	C8-C7-N2-C2
4	B	1303	NAG	C8-C7-N2-C2
4	A	1301	NAG	O7-C7-N2-C2
4	B	1303	NAG	O7-C7-N2-C2
4	A	1302	NAG	O7-C7-N2-C2
4	A	1305	NAG	C8-C7-N2-C2
4	A	1307	NAG	C1-C2-N2-C7
4	C	1303	NAG	C8-C7-N2-C2
4	A	1307	NAG	C3-C2-N2-C7
4	A	1305	NAG	O7-C7-N2-C2
4	C	1303	NAG	O7-C7-N2-C2
4	C	1306	NAG	C8-C7-N2-C2
4	A	1302	NAG	C1-C2-N2-C7
4	A	1303	NAG	C1-C2-N2-C7
4	B	1302	NAG	C1-C2-N2-C7
4	B	1304	NAG	C1-C2-N2-C7
4	C	1304	NAG	C1-C2-N2-C7
4	A	1303	NAG	C3-C2-N2-C7
4	B	1304	NAG	C3-C2-N2-C7
4	B	1308	NAG	C3-C2-N2-C7
4	C	1304	NAG	C3-C2-N2-C7
4	C	1306	NAG	O7-C7-N2-C2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1304	NAG	1	0
4	B	1304	NAG	1	0
4	A	1303	NAG	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-37732. These allow visual inspection of the internal detail of the map and identification of artifacts.

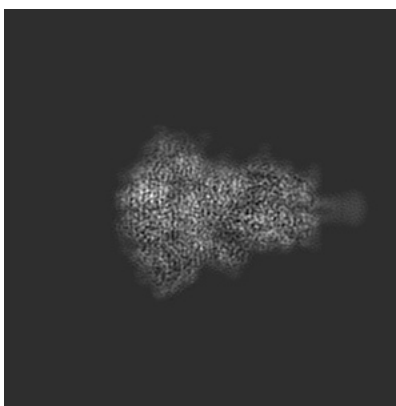
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

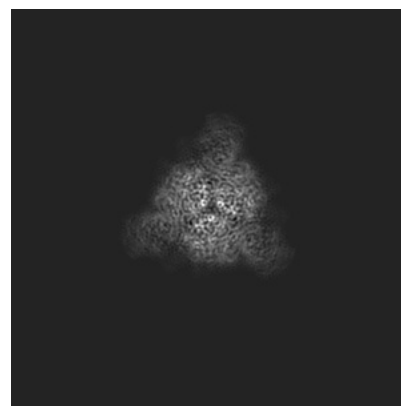
6.1.1 Primary map



X

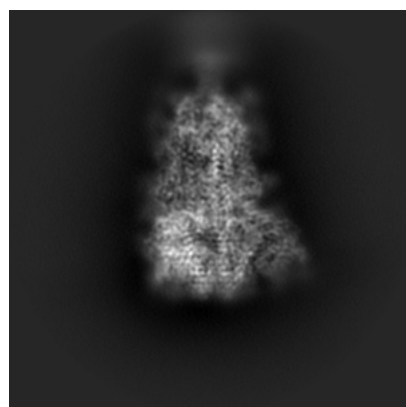


Y

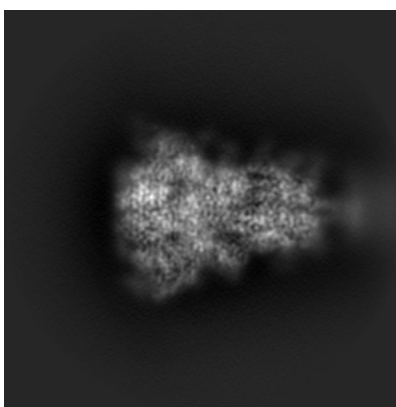


Z

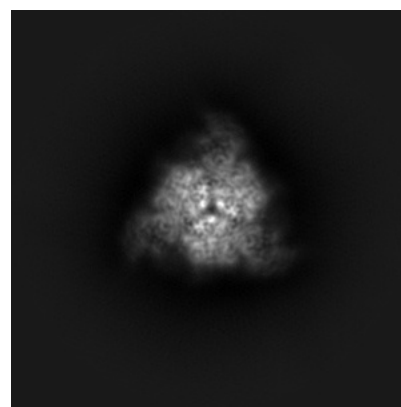
6.1.2 Raw map



X



Y

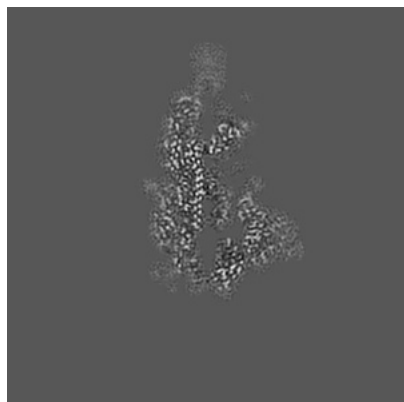


Z

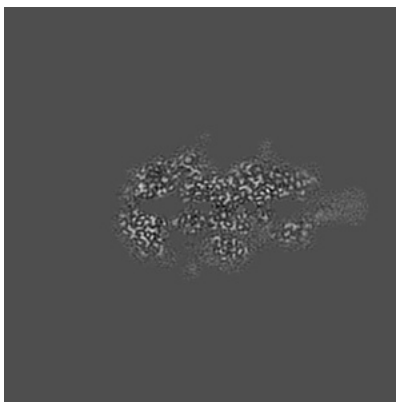
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

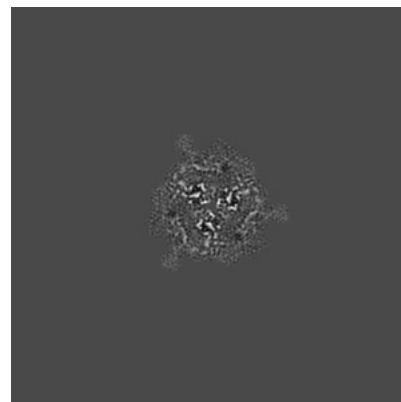
6.2.1 Primary map



X Index: 140

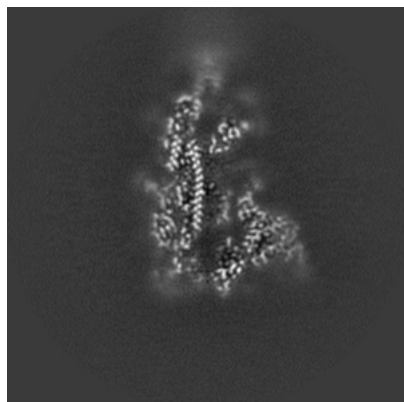


Y Index: 140

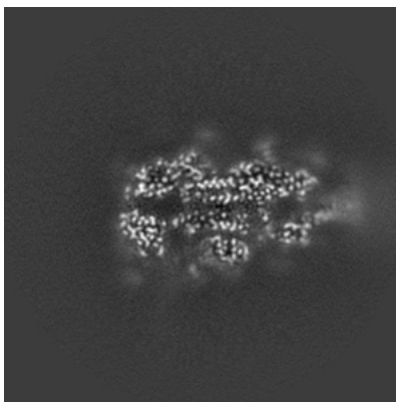


Z Index: 140

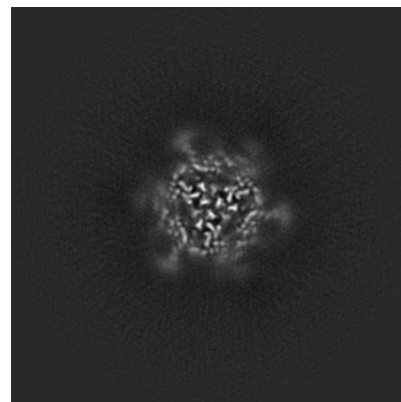
6.2.2 Raw map



X Index: 140



Y Index: 140

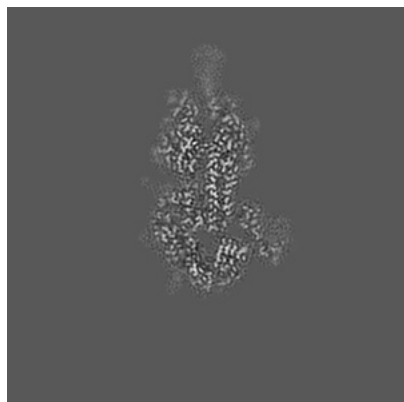


Z Index: 140

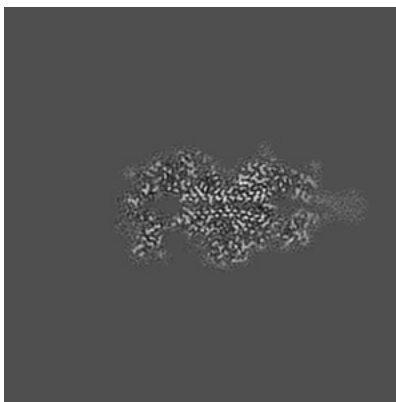
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

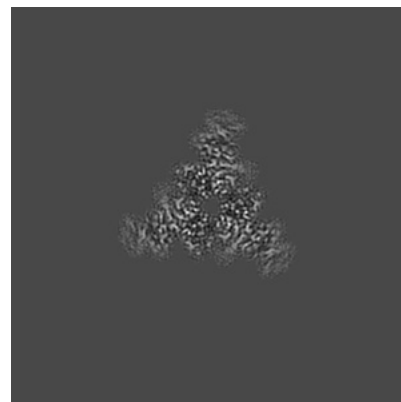
6.3.1 Primary map



X Index: 134

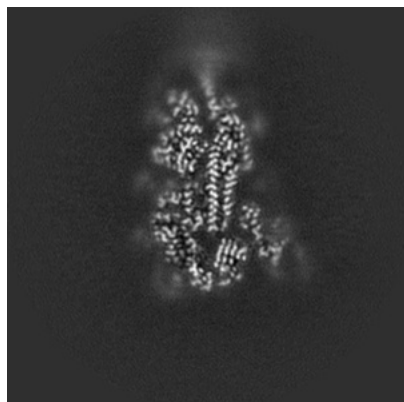


Y Index: 144

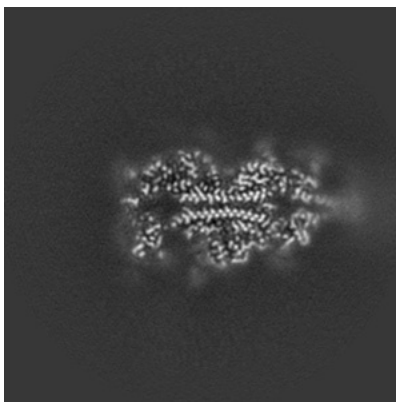


Z Index: 109

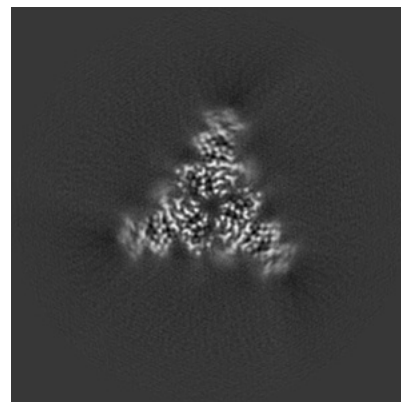
6.3.2 Raw map



X Index: 134



Y Index: 144

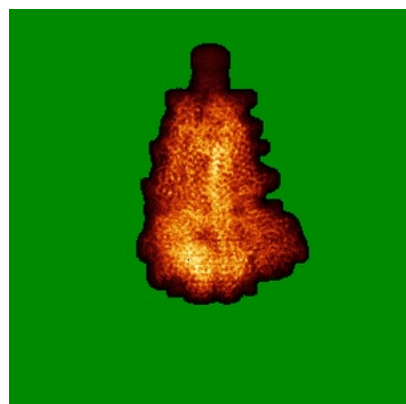


Z Index: 109

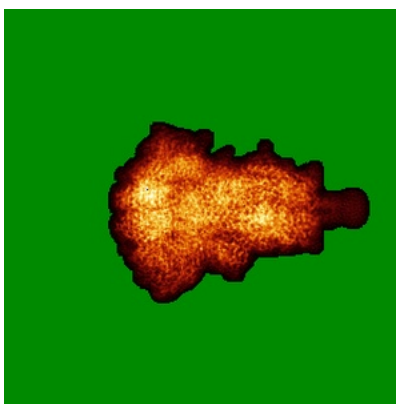
The images above show the largest variance slices of the map in three orthogonal directions.

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

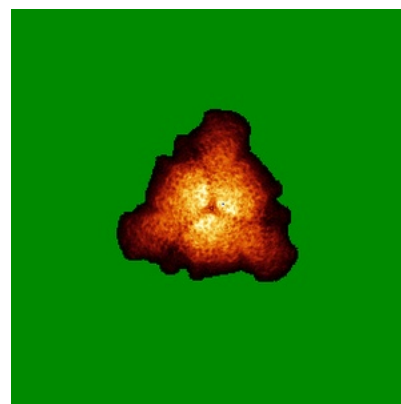
6.4.1 Primary map



X

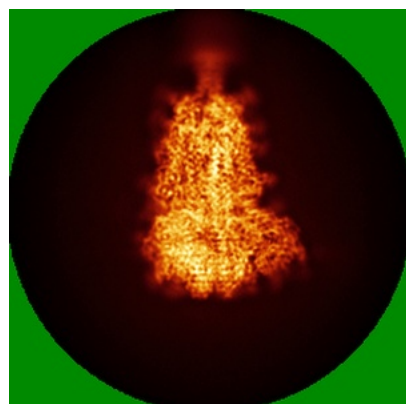


Y

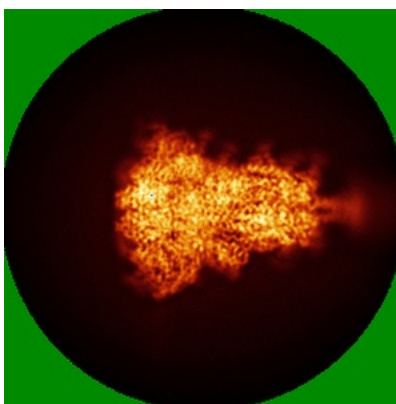


Z

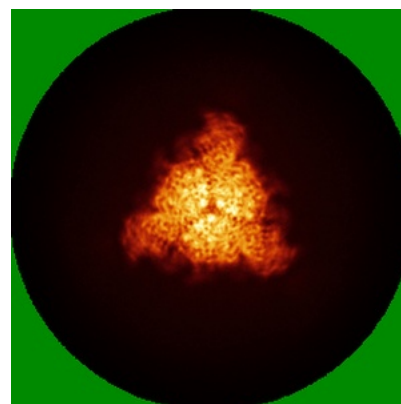
6.4.2 Raw map



X



Y

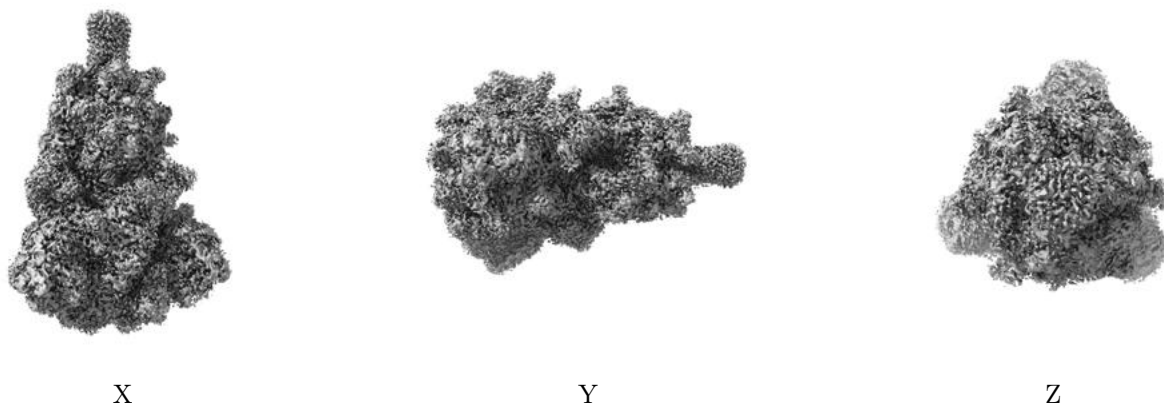


Z

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

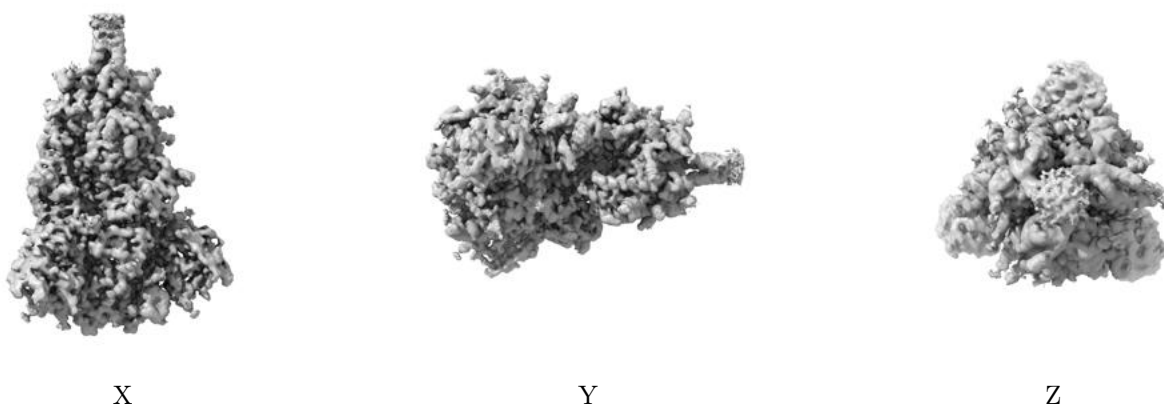
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0115. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

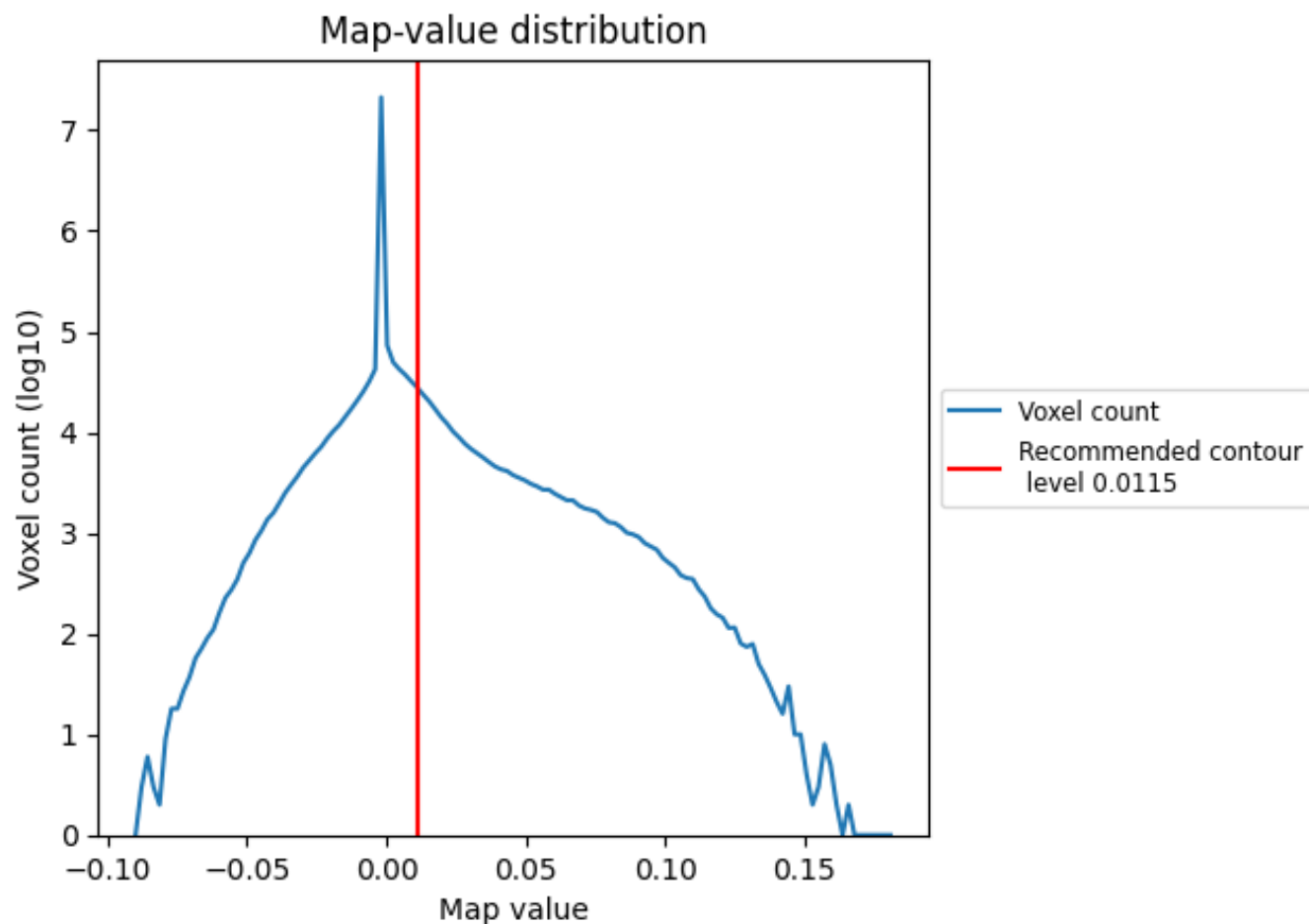
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

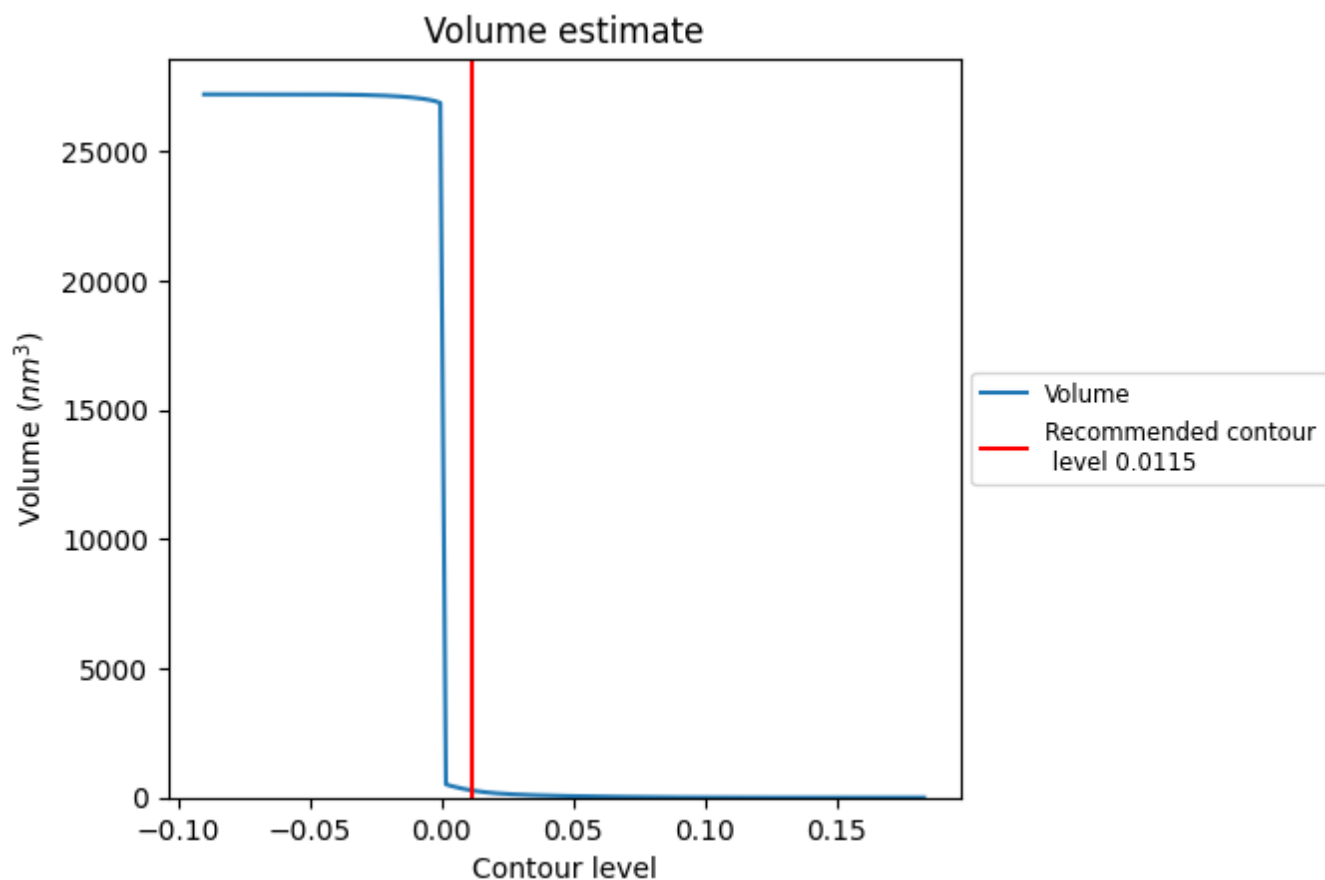
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

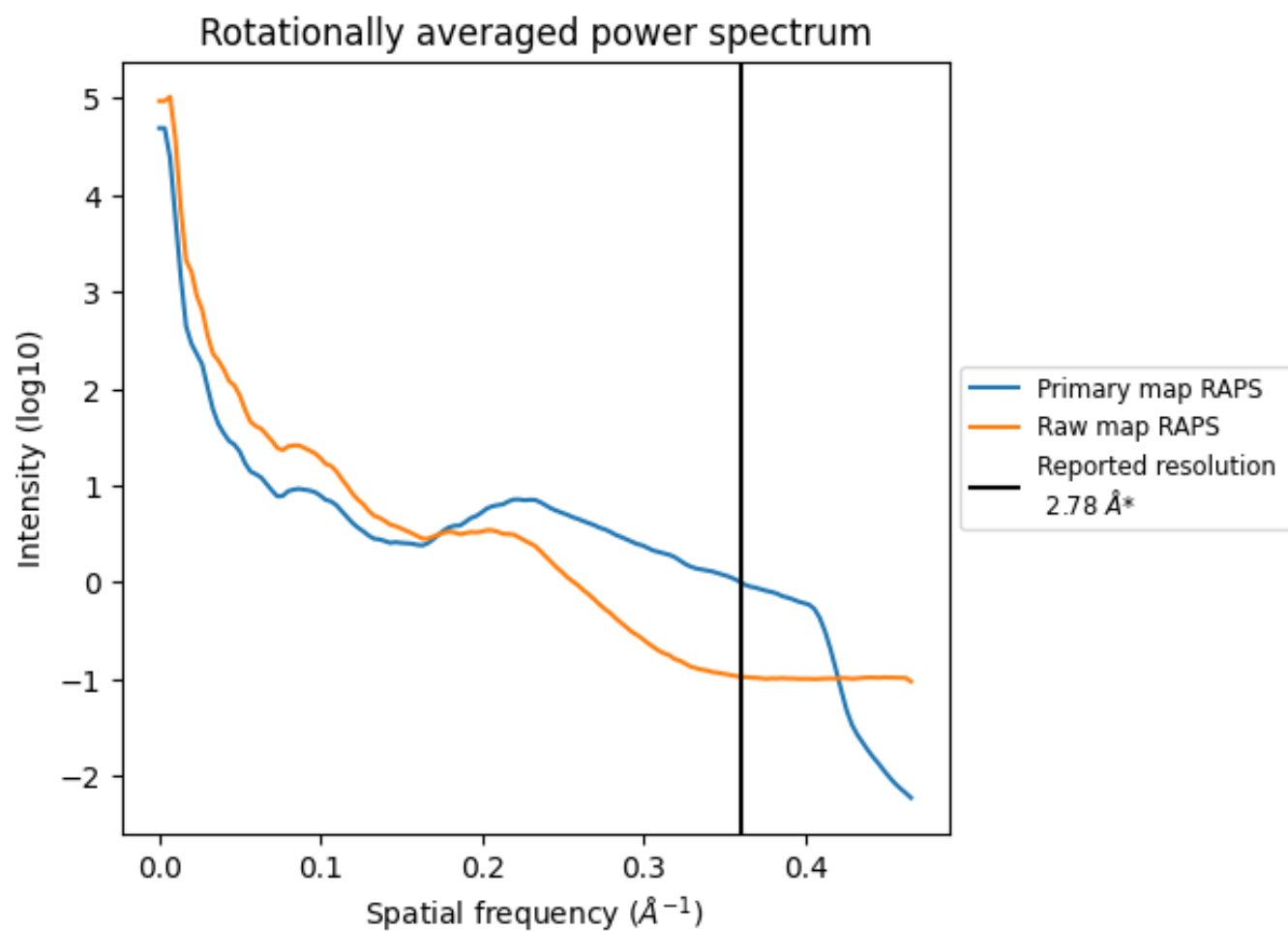
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 279 nm^3 ; this corresponds to an approximate mass of 252 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 ⓘ

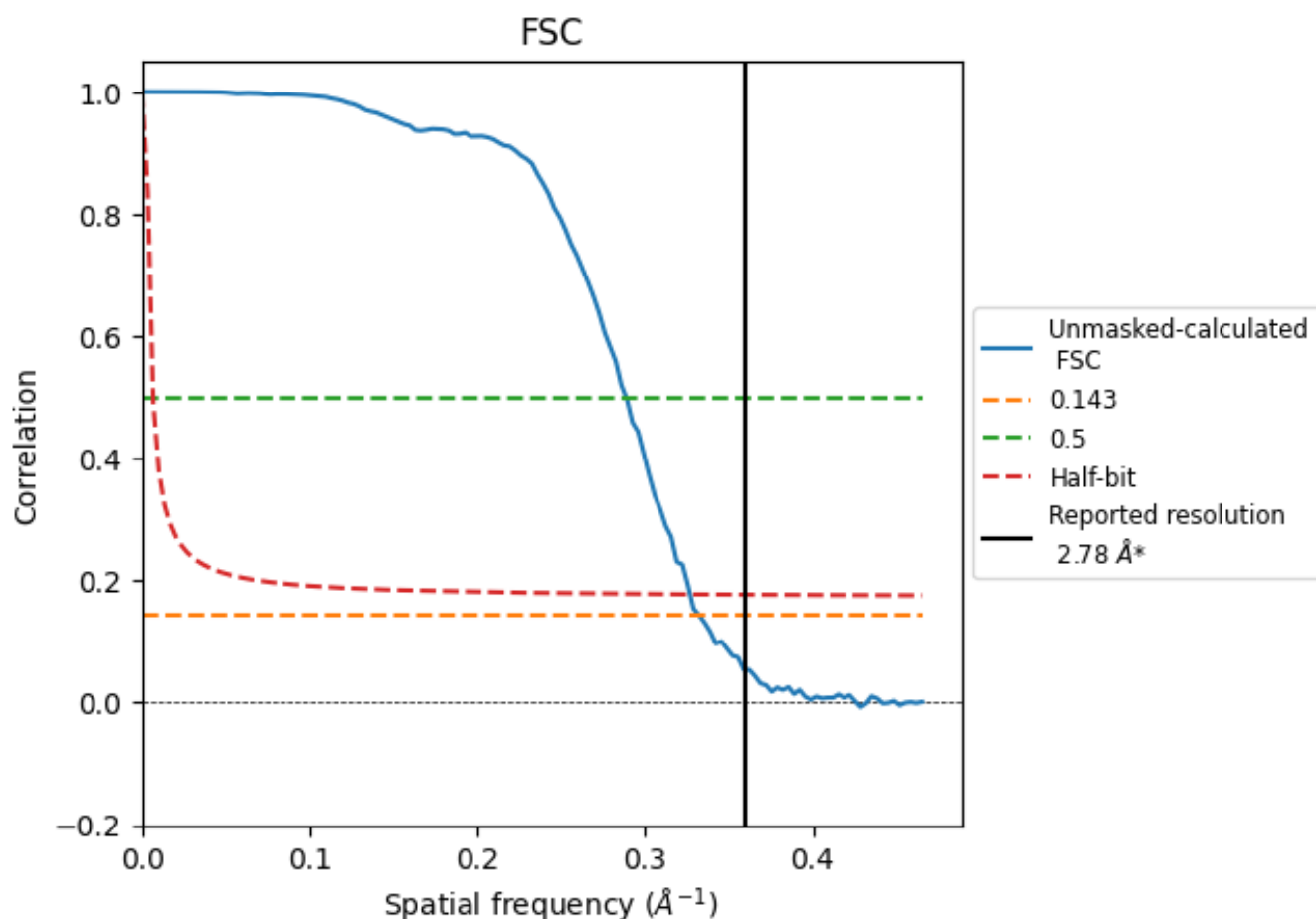


*Reported resolution corresponds to spatial frequency of $0.360 \text{ \AA}^{-1}$

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 0.360 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

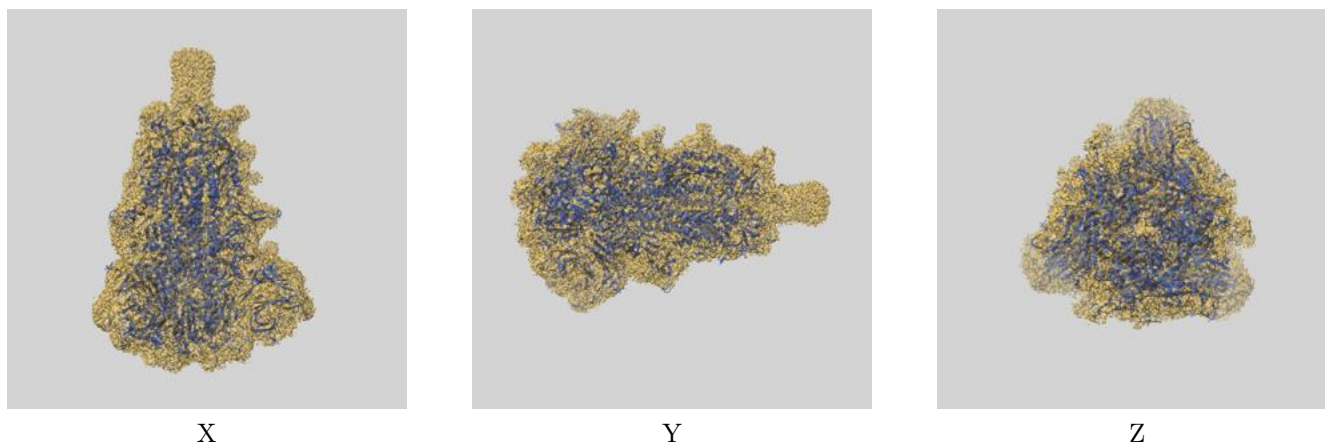
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.78	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.01	3.46	3.06

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

9 Map-model fit [i](#)

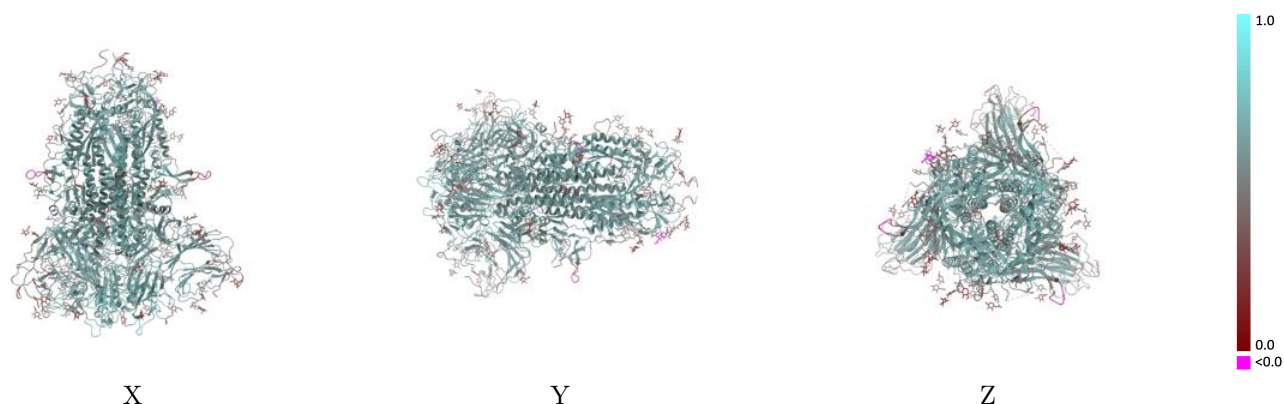
This section contains information regarding the fit between EMDB map EMD-37732 and PDB model 8WQ0. Per-residue inclusion information can be found in section [3](#) on page [15](#).

9.1 Map-model overlay [i](#)



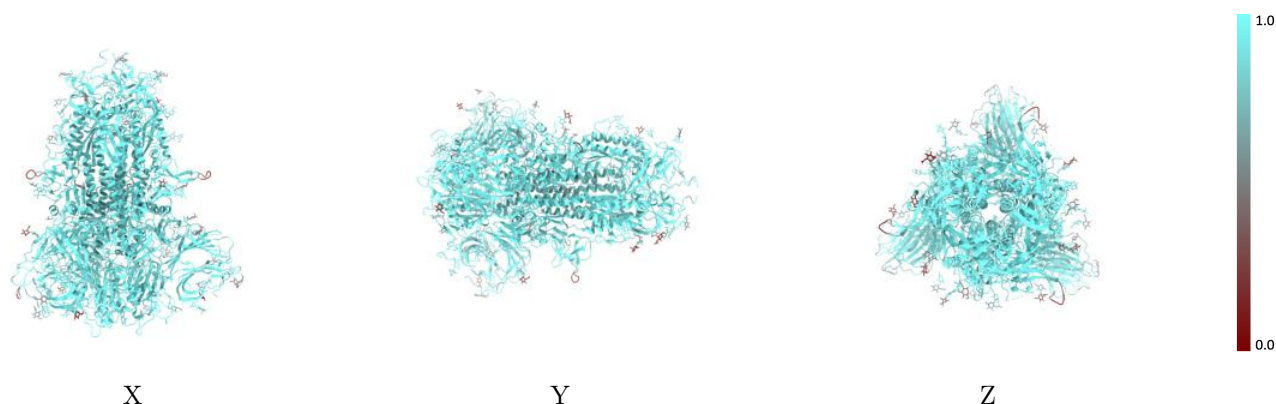
The images above show the 3D surface view of the map at the recommended contour level 0.0115 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



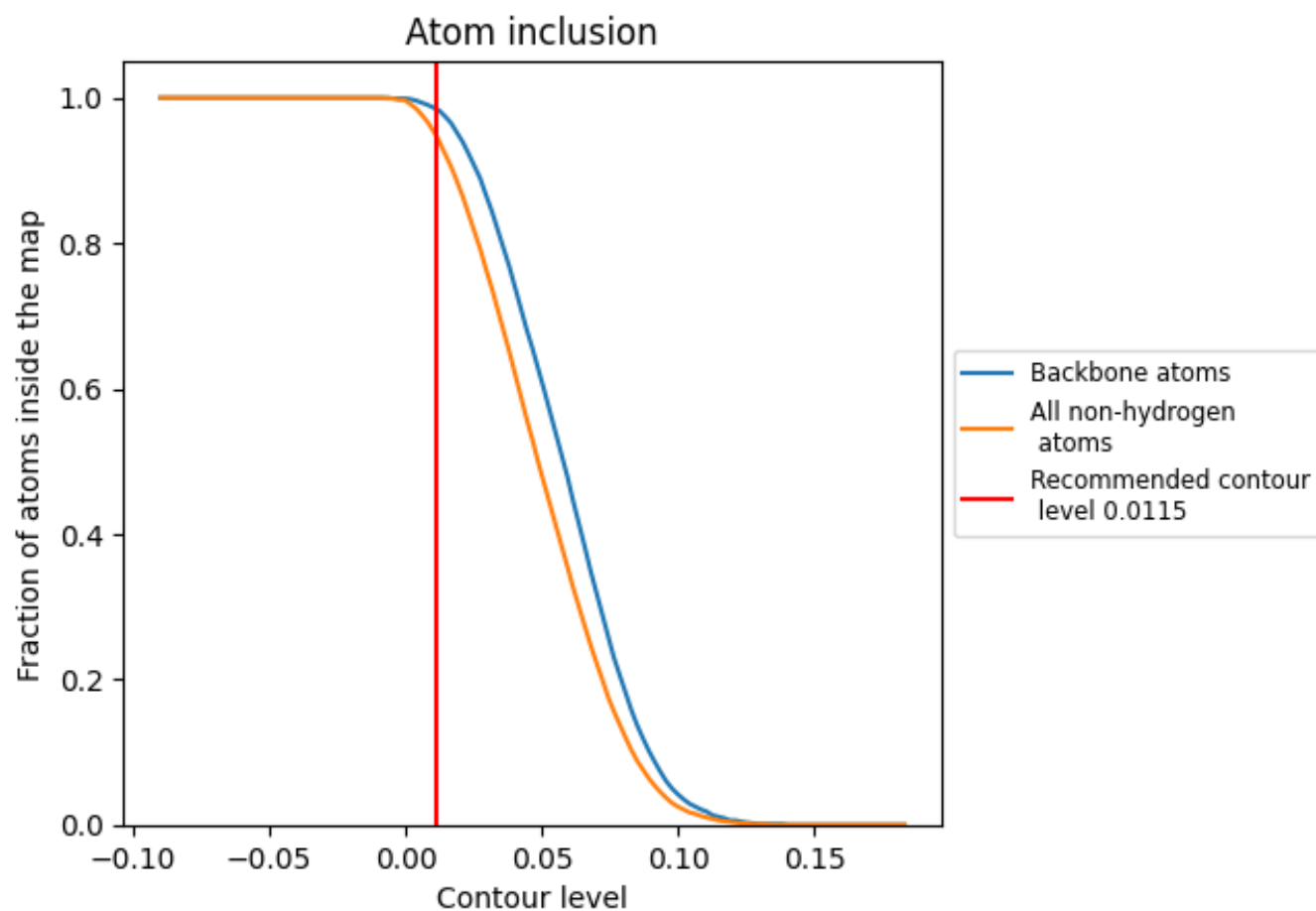
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0115).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 95% 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 (0.0115) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9460	 0.5920
A	 0.9560	 0.6010
B	 0.9540	 0.6020
C	 0.9560	 0.6030
D	 0.5360	 0.2390
E	 0.6410	 0.3910
F	 0.5360	 0.1940
G	 0.7140	 0.2770
H	 0.6790	 0.2520
I	 0.8970	 0.4120
J	 0.8570	 0.3620
K	 0.8930	 0.5200
L	 0.6430	 0.2460
M	 0.8930	 0.3440
N	 0.6430	 0.2830
O	 0.4640	 0.3070
P	 0.7500	 0.4360
Q	 0.6410	 0.3580
R	 0.8570	 0.3640
S	 0.6410	 0.2620
T	 0.6790	 0.2980
U	 0.9290	 0.4860
V	 0.9290	 0.5380
W	 0.7500	 0.3480
X	 0.7500	 0.3300
Y	 0.4290	 0.3110
Z	 0.8210	 0.4510
a	 0.6410	 0.3880
b	 0.9290	 0.5060
c	 0.7500	 0.3750
d	 0.6430	 0.2860
e	 0.2860	 0.1090
f	 0.6410	 0.2430
g	 0.4640	 0.2060
h	 0.7140	 0.2750



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Chain	Atom inclusion	Q-score
i	 0.7140	 0.2710
j	 0.7690	 0.3350