



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 02:27 PM UTC

PDB ID : 8D56 / pdb_00008d56
EMDB ID : EMD-27206
Title : One RBD-up state of SARS-CoV-2 BA.2 variant spike protein
Authors : Zhang, J.; Tang, W.C.; Gao, H.L.; Shi, W.; Peng, H.Q.; Volloch, S.R.; Xiao, T.S.; Chen, B.
Deposited on : 2022-06-04
Resolution : 3.00 Å(reported)
Based on initial model : 7KRR

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

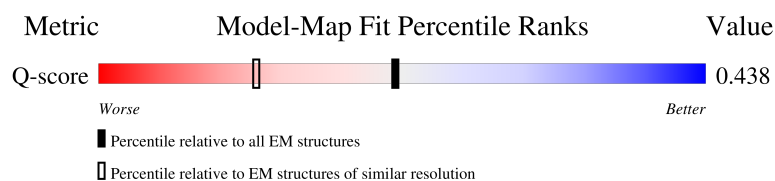
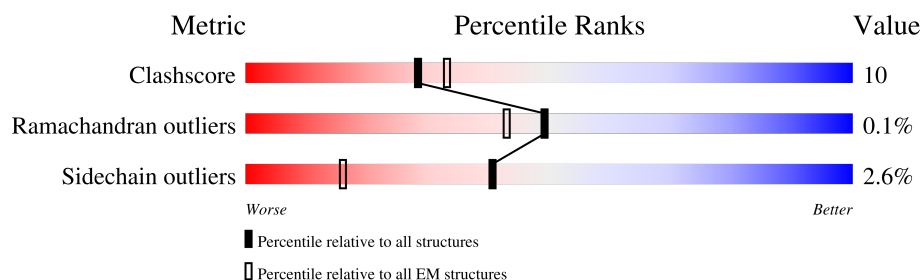
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 3.00 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.



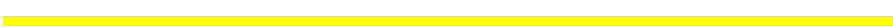
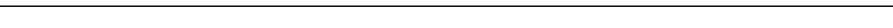
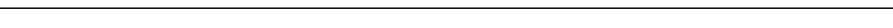
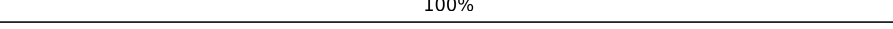
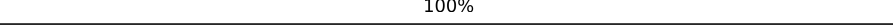
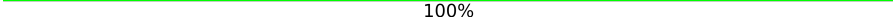
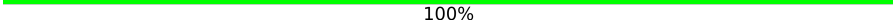
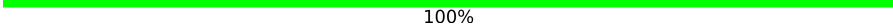
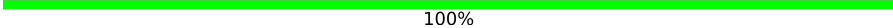
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14081 (2.50 - 3.50)

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	1305	
1	B	1305	
1	C	1305	
2	D	3	

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Mol	Chain	Length	Quality of chain
2	G	3	
2	I	3	
2	J	3	
2	L	3	
2	M	3	
2	N	3	
2	Q	3	
2	S	3	
2	T	3	
2	V	3	
2	W	3	
2	X	3	
2	a	3	
2	c	3	
2	d	3	
2	f	3	
2	g	3	
3	E	2	
3	F	2	
3	H	2	
3	O	2	
3	P	2	
3	R	2	
3	Y	2	
3	Z	2	

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Mol	Chain	Length	Quality of chain
3	b	2	50% 50%
4	K	3	33% 67%
4	U	3	33% 67%
4	e	3	33% 67%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 27674 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	1110	Total	C	N	O	S	0	0
			8715	5575	1455	1646	39		
1	B	1124	Total	C	N	O	S	0	0
			8820	5640	1473	1667	40		
1	C	1119	Total	C	N	O	S	0	0
			8777	5613	1465	1659	40		

There are 198 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	ILE	THR	variant	UNP P0DTC2
A	?	-	LEU	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	27	SER	ALA	variant	UNP P0DTC2
A	142	ASP	GLY	variant	UNP P0DTC2
A	213	GLY	VAL	variant	UNP P0DTC2
A	339	ASP	GLY	variant	UNP P0DTC2
A	371	PHE	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	376	ALA	THR	variant	UNP P0DTC2
A	405	ASN	ASP	variant	UNP P0DTC2
A	408	SER	ARG	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	484	ALA	GLU	variant	UNP P0DTC2
A	493	ARG	GLN	variant	UNP P0DTC2
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	655	TYR	HIS	variant	UNP P0DTC2
A	679	LYS	ASN	variant	UNP P0DTC2
A	681	HIS	PRO	variant	UNP P0DTC2
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	954	HIS	GLN	variant	UNP P0DTC2
A	969	LYS	ASN	variant	UNP P0DTC2
A	1274	SER	-	expression tag	UNP P0DTC2
A	1275	GLY	-	expression tag	UNP P0DTC2
A	1276	GLY	-	expression tag	UNP P0DTC2
A	1277	GLY	-	expression tag	UNP P0DTC2
A	1278	SER	-	expression tag	UNP P0DTC2
A	1279	ALA	-	expression tag	UNP P0DTC2
A	1280	TRP	-	expression tag	UNP P0DTC2
A	1281	SER	-	expression tag	UNP P0DTC2
A	1282	HIS	-	expression tag	UNP P0DTC2
A	1283	PRO	-	expression tag	UNP P0DTC2
A	1284	GLN	-	expression tag	UNP P0DTC2
A	1285	PHE	-	expression tag	UNP P0DTC2
A	1286	GLU	-	expression tag	UNP P0DTC2
A	1287	LYS	-	expression tag	UNP P0DTC2
A	1288	GLY	-	expression tag	UNP P0DTC2
A	1289	GLY	-	expression tag	UNP P0DTC2
A	1290	GLY	-	expression tag	UNP P0DTC2
A	1291	SER	-	expression tag	UNP P0DTC2
A	1292	GLY	-	expression tag	UNP P0DTC2
A	1293	GLY	-	expression tag	UNP P0DTC2
A	1294	GLY	-	expression tag	UNP P0DTC2
A	1295	SER	-	expression tag	UNP P0DTC2
A	1296	GLY	-	expression tag	UNP P0DTC2
A	1297	GLY	-	expression tag	UNP P0DTC2
A	1298	SER	-	expression tag	UNP P0DTC2
A	1299	SER	-	expression tag	UNP P0DTC2
A	1300	ALA	-	expression tag	UNP P0DTC2
A	1301	TRP	-	expression tag	UNP P0DTC2
A	1302	SER	-	expression tag	UNP P0DTC2
A	1303	HIS	-	expression tag	UNP P0DTC2
A	1304	PRO	-	expression tag	UNP P0DTC2
A	1305	GLN	-	expression tag	UNP P0DTC2
A	1306	PHE	-	expression tag	UNP P0DTC2
A	1307	GLU	-	expression tag	UNP P0DTC2
A	1308	LYS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	19	ILE	THR	variant	UNP P0DTC2
B	?	-	LEU	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	27	SER	ALA	variant	UNP P0DTC2
B	142	ASP	GLY	variant	UNP P0DTC2
B	213	GLY	VAL	variant	UNP P0DTC2
B	339	ASP	GLY	variant	UNP P0DTC2
B	371	PHE	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	376	ALA	THR	variant	UNP P0DTC2
B	405	ASN	ASP	variant	UNP P0DTC2
B	408	SER	ARG	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	477	ASN	SER	variant	UNP P0DTC2
B	478	LYS	THR	variant	UNP P0DTC2
B	484	ALA	GLU	variant	UNP P0DTC2
B	493	ARG	GLN	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	505	HIS	TYR	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	679	LYS	ASN	variant	UNP P0DTC2
B	681	HIS	PRO	variant	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	954	HIS	GLN	variant	UNP P0DTC2
B	969	LYS	ASN	variant	UNP P0DTC2
B	1274	SER	-	expression tag	UNP P0DTC2
B	1275	GLY	-	expression tag	UNP P0DTC2
B	1276	GLY	-	expression tag	UNP P0DTC2
B	1277	GLY	-	expression tag	UNP P0DTC2
B	1278	SER	-	expression tag	UNP P0DTC2
B	1279	ALA	-	expression tag	UNP P0DTC2
B	1280	TRP	-	expression tag	UNP P0DTC2
B	1281	SER	-	expression tag	UNP P0DTC2
B	1282	HIS	-	expression tag	UNP P0DTC2
B	1283	PRO	-	expression tag	UNP P0DTC2
B	1284	GLN	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1285	PHE	-	expression tag	UNP P0DTC2
B	1286	GLU	-	expression tag	UNP P0DTC2
B	1287	LYS	-	expression tag	UNP P0DTC2
B	1288	GLY	-	expression tag	UNP P0DTC2
B	1289	GLY	-	expression tag	UNP P0DTC2
B	1290	GLY	-	expression tag	UNP P0DTC2
B	1291	SER	-	expression tag	UNP P0DTC2
B	1292	GLY	-	expression tag	UNP P0DTC2
B	1293	GLY	-	expression tag	UNP P0DTC2
B	1294	GLY	-	expression tag	UNP P0DTC2
B	1295	SER	-	expression tag	UNP P0DTC2
B	1296	GLY	-	expression tag	UNP P0DTC2
B	1297	GLY	-	expression tag	UNP P0DTC2
B	1298	SER	-	expression tag	UNP P0DTC2
B	1299	SER	-	expression tag	UNP P0DTC2
B	1300	ALA	-	expression tag	UNP P0DTC2
B	1301	TRP	-	expression tag	UNP P0DTC2
B	1302	SER	-	expression tag	UNP P0DTC2
B	1303	HIS	-	expression tag	UNP P0DTC2
B	1304	PRO	-	expression tag	UNP P0DTC2
B	1305	GLN	-	expression tag	UNP P0DTC2
B	1306	PHE	-	expression tag	UNP P0DTC2
B	1307	GLU	-	expression tag	UNP P0DTC2
B	1308	LYS	-	expression tag	UNP P0DTC2
C	19	ILE	THR	variant	UNP P0DTC2
C	?	-	LEU	deletion	UNP P0DTC2
C	?	-	PRO	deletion	UNP P0DTC2
C	?	-	PRO	deletion	UNP P0DTC2
C	27	SER	ALA	variant	UNP P0DTC2
C	142	ASP	GLY	variant	UNP P0DTC2
C	213	GLY	VAL	variant	UNP P0DTC2
C	339	ASP	GLY	variant	UNP P0DTC2
C	371	PHE	SER	variant	UNP P0DTC2
C	373	PRO	SER	variant	UNP P0DTC2
C	375	PHE	SER	variant	UNP P0DTC2
C	376	ALA	THR	variant	UNP P0DTC2
C	405	ASN	ASP	variant	UNP P0DTC2
C	408	SER	ARG	variant	UNP P0DTC2
C	417	ASN	LYS	variant	UNP P0DTC2
C	440	LYS	ASN	variant	UNP P0DTC2
C	477	ASN	SER	variant	UNP P0DTC2
C	478	LYS	THR	variant	UNP P0DTC2

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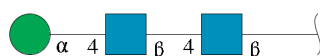
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C	484	ALA	GLU	variant	UNP P0DTC2
C	493	ARG	GLN	variant	UNP P0DTC2
C	498	ARG	GLN	variant	UNP P0DTC2
C	501	TYR	ASN	variant	UNP P0DTC2
C	505	HIS	TYR	variant	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	655	TYR	HIS	variant	UNP P0DTC2
C	679	LYS	ASN	variant	UNP P0DTC2
C	681	HIS	PRO	variant	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2
C	954	HIS	GLN	variant	UNP P0DTC2
C	969	LYS	ASN	variant	UNP P0DTC2
C	1274	SER	-	expression tag	UNP P0DTC2
C	1275	GLY	-	expression tag	UNP P0DTC2
C	1276	GLY	-	expression tag	UNP P0DTC2
C	1277	GLY	-	expression tag	UNP P0DTC2
C	1278	SER	-	expression tag	UNP P0DTC2
C	1279	ALA	-	expression tag	UNP P0DTC2
C	1280	TRP	-	expression tag	UNP P0DTC2
C	1281	SER	-	expression tag	UNP P0DTC2
C	1282	HIS	-	expression tag	UNP P0DTC2
C	1283	PRO	-	expression tag	UNP P0DTC2
C	1284	GLN	-	expression tag	UNP P0DTC2
C	1285	PHE	-	expression tag	UNP P0DTC2
C	1286	GLU	-	expression tag	UNP P0DTC2
C	1287	LYS	-	expression tag	UNP P0DTC2
C	1288	GLY	-	expression tag	UNP P0DTC2
C	1289	GLY	-	expression tag	UNP P0DTC2
C	1290	GLY	-	expression tag	UNP P0DTC2
C	1291	SER	-	expression tag	UNP P0DTC2
C	1292	GLY	-	expression tag	UNP P0DTC2
C	1293	GLY	-	expression tag	UNP P0DTC2
C	1294	GLY	-	expression tag	UNP P0DTC2
C	1295	SER	-	expression tag	UNP P0DTC2
C	1296	GLY	-	expression tag	UNP P0DTC2
C	1297	GLY	-	expression tag	UNP P0DTC2
C	1298	SER	-	expression tag	UNP P0DTC2
C	1299	SER	-	expression tag	UNP P0DTC2
C	1300	ALA	-	expression tag	UNP P0DTC2
C	1301	TRP	-	expression tag	UNP P0DTC2
C	1302	SER	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1303	HIS	-	expression tag	UNP P0DTC2
C	1304	PRO	-	expression tag	UNP P0DTC2
C	1305	GLN	-	expression tag	UNP P0DTC2
C	1306	PHE	-	expression tag	UNP P0DTC2
C	1307	GLU	-	expression tag	UNP P0DTC2
C	1308	LYS	-	expression tag	UNP P0DTC2

- Molecule 2 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.



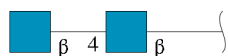
Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	3	Total	C	N	O	0	0
			39	22	2	15		
2	G	3	Total	C	N	O	0	0
			39	22	2	15		
2	I	3	Total	C	N	O	0	0
			39	22	2	15		
2	J	3	Total	C	N	O	0	0
			39	22	2	15		
2	L	3	Total	C	N	O	0	0
			39	22	2	15		
2	M	3	Total	C	N	O	0	0
			39	22	2	15		
2	N	3	Total	C	N	O	0	0
			39	22	2	15		
2	Q	3	Total	C	N	O	0	0
			39	22	2	15		
2	S	3	Total	C	N	O	0	0
			39	22	2	15		
2	T	3	Total	C	N	O	0	0
			39	22	2	15		
2	V	3	Total	C	N	O	0	0
			39	22	2	15		
2	W	3	Total	C	N	O	0	0
			39	22	2	15		
2	X	3	Total	C	N	O	0	0
			39	22	2	15		

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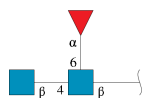
Mol	Chain	Residues	Atoms				AltConf	Trace
2	a	3	Total	C	N	O	0	0
			39	22	2	15		
2	c	3	Total	C	N	O	0	0
			39	22	2	15		
2	d	3	Total	C	N	O	0	0
			39	22	2	15		
2	f	3	Total	C	N	O	0	0
			39	22	2	15		
2	g	3	Total	C	N	O	0	0
			39	22	2	15		

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



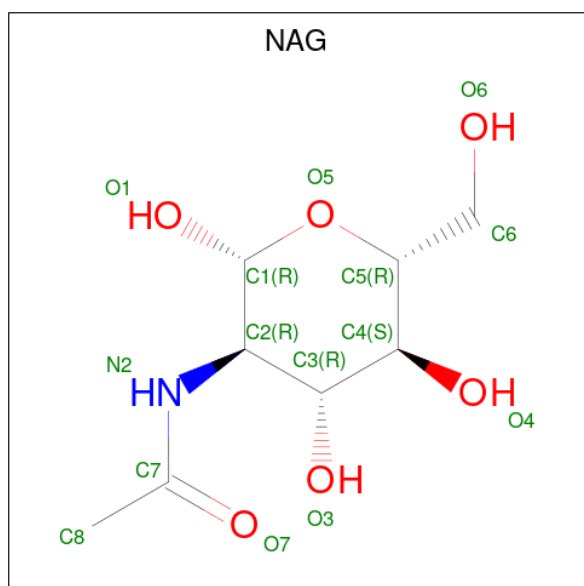
Mol	Chain	Residues	Atoms				AltConf	Trace
3	E	2	Total	C	N	O	0	0
			28	16	2	10		
3	F	2	Total	C	N	O	0	0
			28	16	2	10		
3	H	2	Total	C	N	O	0	0
			28	16	2	10		
3	O	2	Total	C	N	O	0	0
			28	16	2	10		
3	P	2	Total	C	N	O	0	0
			28	16	2	10		
3	R	2	Total	C	N	O	0	0
			28	16	2	10		
3	Y	2	Total	C	N	O	0	0
			28	16	2	10		
3	Z	2	Total	C	N	O	0	0
			28	16	2	10		
3	b	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 4 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
4	K	3	Total	C	N	O	0	0
			38	22	2	14		
4	U	3	Total	C	N	O	0	0
			38	22	2	14		
4	e	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).



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

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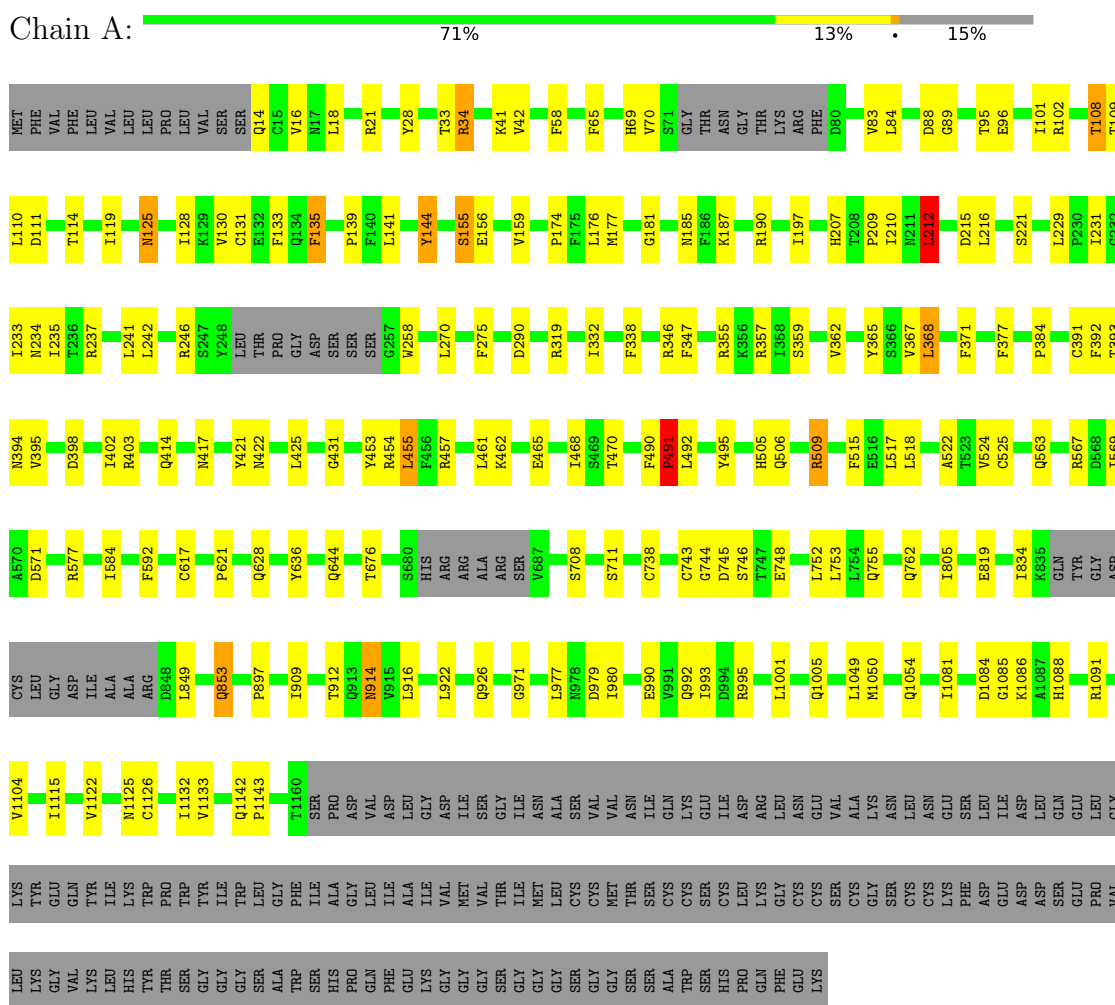
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Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C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

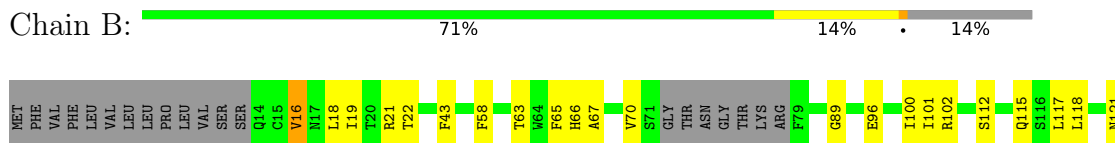
3 Residue-property plots

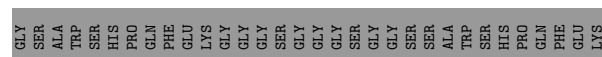
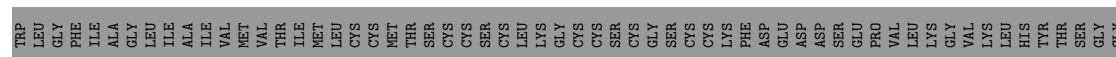
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: Spike glycoprotein



• Molecule 1: Spike glycoprotein





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

Chain D: 100%



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

Chain G: 33% 67%



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

Chain I: 67% 33%



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

Chain J: 67% 33%



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

Chain L: 67% 33%



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

Chain M:  67% 33%



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

Chain N:  100%



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

Chain Q:  67% 33%



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

Chain S:  67% 33%



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

Chain T:  67% 33%



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

Chain V:  67% 33%

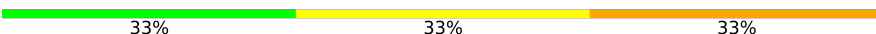


- Molecule 2: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain W:  67% 33%



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

Chain X:  33% 33% 33%



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

Chain a:  67% 33%



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

Chain c:  67% 33%



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

Chain d:  33% 67%



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

Chain f:  33% 67%

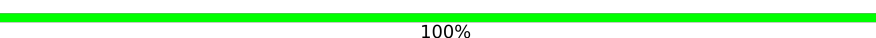


- Molecule 2: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain g:  67% 33%



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

Chain E:  100%



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

Chain F:  100%



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

Chain H:  50% 50%



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

Chain O:  100%



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

Chain P:  100%



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

Chain R:  100%



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

Chain Y:  100%



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

Chain Z:  100%

MAG1
MAG2

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

Chain b:  50% 50%

MAG1
MAG2

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

Chain K:  33%  67%

MAG1
MAG2
FUC3

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

Chain U:  33%  67%

MAG1
MAG2
FUC3

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

Chain e:  33%  67%

MAG1
MAG2
FUC3

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	193655	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 ($e^-/\text{\AA}^2$)	53.853	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.556	Depositor
Minimum map value	-0.187	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.012	Depositor
Recommended contour level	0.03	Depositor
Map size (Å)	398.4, 398.4, 398.4	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	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: MAN, NAG, FUC

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.72	0/8921	1.19	12/12133 (0.1%)
1	B	0.74	1/9029 (0.0%)	1.20	9/12280 (0.1%)
1	C	0.75	0/8984	1.24	19/12219 (0.2%)
All	All	0.74	1/26934 (0.0%)	1.21	40/36632 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	133	PHE	N-CA	-5.17	1.39	1.45

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	172	SER	N-CA-C	8.55	119.22	108.45
1	B	216	LEU	CB-CA-C	7.35	116.42	111.20
1	C	246	ARG	N-CA-C	-7.29	102.33	112.25
1	A	125	ASN	N-CA-C	7.25	120.22	108.32
1	C	125	ASN	N-CA-C	7.19	117.51	108.45
1	C	1048	HIS	N-CA-C	7.08	119.99	109.59
1	B	852	ALA	N-CA-C	-6.91	103.83	111.36
1	A	1091	ARG	N-CA-C	-6.72	103.57	111.03
1	A	979	ASP	N-CA-C	-6.43	103.89	111.03
1	B	490	PHE	N-CA-C	-6.41	97.85	109.44
1	A	135	PHE	N-CA-C	-6.37	101.06	110.48
1	C	334	ASN	N-CA-C	6.23	120.25	111.52
1	C	493	ARG	N-CA-C	6.19	119.02	110.35
1	A	628	GLN	N-CA-C	6.02	119.19	110.28
1	C	170	TYR	N-CA-C	6.01	117.98	108.79
1	A	131	CYS	N-CA-C	6.00	118.09	109.14
1	C	677	GLN	N-CA-C	5.99	117.68	111.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	19	ILE	N-CA-C	5.90	116.42	108.17
1	A	455	LEU	N-CA-C	-5.65	104.98	112.94
1	C	158	ARG	CA-C-N	5.64	128.19	120.46
1	C	158	ARG	C-N-CA	5.64	128.19	120.46
1	B	214	ARG	N-CA-C	5.55	117.53	108.55
1	C	236	THR	N-CA-C	5.49	120.09	113.17
1	A	181	GLY	N-CA-C	-5.44	102.76	112.22
1	A	108	THR	N-CA-C	5.43	117.28	111.36
1	B	577	ARG	N-CA-C	-5.39	100.39	109.07
1	A	212	LEU	N-CA-C	-5.37	106.68	113.23
1	B	89	GLY	N-CA-C	5.31	117.00	111.95
1	B	622	VAL	N-CA-C	-5.27	104.38	111.44
1	B	425	LEU	CB-CA-C	5.23	116.68	108.63
1	C	516	GLU	N-CA-C	5.22	119.14	112.87
1	C	229	LEU	CA-C-N	5.18	126.32	119.84
1	C	229	LEU	C-N-CA	5.18	126.32	119.84
1	A	853	GLN	N-CA-C	-5.17	105.34	110.97
1	A	34	ARG	N-CA-C	5.09	117.25	110.53
1	B	620	VAL	N-CA-CB	5.08	113.70	110.50
1	C	247	SER	CA-C-N	-5.08	114.84	122.81
1	C	247	SER	C-N-CA	-5.08	114.84	122.81
1	C	833	PHE	CA-C-N	5.02	127.42	120.49
1	C	833	PHE	C-N-CA	5.02	127.42	120.49

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	8715	0	8502	156	0
1	B	8820	0	8596	217	0
1	C	8777	0	8562	165	0
2	D	39	0	34	5	0
2	G	39	0	34	1	0
2	I	39	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	J	39	0	34	0	0
2	L	39	0	34	0	0
2	M	39	0	34	0	0
2	N	39	0	34	2	0
2	Q	39	0	34	0	0
2	S	39	0	34	0	0
2	T	39	0	34	0	0
2	V	39	0	34	0	0
2	W	39	0	34	0	0
2	X	39	0	34	1	0
2	a	39	0	34	0	0
2	c	39	0	34	0	0
2	d	39	0	34	1	0
2	f	39	0	34	1	0
2	g	39	0	34	0	0
3	E	28	0	25	0	0
3	F	28	0	25	0	0
3	H	28	0	25	2	0
3	O	28	0	25	0	0
3	P	28	0	25	0	0
3	R	28	0	25	0	0
3	Y	28	0	25	0	0
3	Z	28	0	25	0	0
3	b	28	0	25	0	0
4	K	38	0	34	0	0
4	U	38	0	34	0	0
4	e	38	0	34	0	0
5	A	98	0	91	2	0
5	B	98	0	91	2	0
5	C	98	0	91	0	0
All	All	27674	0	26872	518	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 10.

All (518) 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:454:ARG:NH2	1:B:491:PRO:HG2	1.39	1.36
1:B:452:LEU:CD2	1:B:492:LEU:HD22	1.71	1.18
1:B:454:ARG:CZ	1:B:491:PRO:HG2	1.84	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:140:PHE:HE1	1:C:244:LEU:HB2	1.15	1.03
1:B:493:ARG:HH12	1:C:378:LYS:HD3	1.19	1.03
1:A:33:THR:HG22	1:A:58:PHE:CE2	1.94	1.01
1:C:310:LYS:HA	1:C:600:PRO:HA	1.44	1.00
1:B:454:ARG:NH1	1:B:491:PRO:HG3	1.75	0.99
1:B:493:ARG:HH12	1:C:378:LYS:CD	1.74	0.99
1:B:452:LEU:HD21	1:B:492:LEU:HD22	1.40	0.98
1:A:393:THR:HA	1:A:522:ALA:HA	1.44	0.96
1:B:454:ARG:NH2	1:B:491:PRO:CG	2.31	0.94
1:A:33:THR:HG22	1:A:58:PHE:HE2	1.29	0.93
1:B:454:ARG:CZ	1:B:491:PRO:CG	2.45	0.93
1:B:493:ARG:HH22	1:C:378:LYS:HD2	1.35	0.92
1:B:454:ARG:HH22	1:B:491:PRO:HG2	1.34	0.91
1:C:515:PHE:O	1:C:517:LEU:HD12	1.72	0.89
1:B:456:PHE:O	1:B:491:PRO:HG3	1.72	0.88
1:B:452:LEU:HD11	1:B:493:ARG:N	1.89	0.87
1:C:140:PHE:CE1	1:C:244:LEU:HB2	2.08	0.86
1:C:515:PHE:O	1:C:517:LEU:CD1	2.26	0.83
1:A:490:PHE:HB2	1:A:491:PRO:HD2	1.59	0.82
1:B:452:LEU:HD21	1:B:492:LEU:CD2	2.10	0.81
1:B:452:LEU:HD23	1:B:492:LEU:HD22	1.62	0.81
1:B:454:ARG:NH1	1:B:491:PRO:CG	2.43	0.81
1:B:456:PHE:HB2	1:B:491:PRO:HB3	1.63	0.79
1:C:347:PHE:CD2	1:C:399:SER:HB2	2.19	0.78
1:B:454:ARG:HH12	1:B:491:PRO:HG3	1.49	0.77
1:A:834:ILE:O	1:A:834:ILE:HD12	1.85	0.76
1:C:1054:GLN:HE21	1:C:1054:GLN:HA	1.50	0.76
1:A:1054:GLN:HA	1:A:1054:GLN:HE21	1.51	0.76
1:A:28:TYR:CE2	5:A:1401:NAG:H5	2.21	0.75
1:B:457:ARG:HH12	1:B:461:LEU:HD12	1.52	0.75
1:B:112:SER:O	1:B:132:GLU:OE2	2.05	0.75
1:B:493:ARG:NH2	1:C:378:LYS:HD2	2.02	0.74
1:C:98:SER:O	1:C:102:ARG:NH2	2.20	0.74
1:A:417:ASN:HB2	1:A:455:LEU:HD22	1.68	0.74
1:C:677:GLN:CD	1:C:677:GLN:H	1.95	0.73
1:C:347:PHE:HD2	1:C:399:SER:HB2	1.54	0.73
1:C:520:ALA:HB3	1:C:521:PRO:HD3	1.69	0.73
1:B:834:ILE:HG21	3:H:1:NAG:H81	1.70	0.73
1:C:453:TYR:OH	1:C:493:ARG:NH1	2.22	0.73
1:B:456:PHE:H	1:B:491:PRO:HB3	1.54	0.73
1:C:120:VAL:HG12	1:C:129:LYS:NZ	2.04	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1085:GLY:O	1:A:1125:ASN:HA	1.90	0.72
1:B:121:ASN:HD21	1:B:174:PRO:HB3	1.54	0.72
1:B:449:TYR:HA	1:B:494:SER:OG	1.90	0.72
1:A:377:PHE:HE2	1:A:384:PRO:HB3	1.53	0.71
1:B:452:LEU:HD11	1:B:493:ARG:H	1.54	0.70
1:B:493:ARG:NH1	1:C:378:LYS:CD	2.50	0.70
1:A:971:GLY:HA3	1:A:995:ARG:HH11	1.56	0.70
1:B:457:ARG:NH1	1:B:461:LEU:HD12	2.07	0.70
1:B:577:ARG:NH1	1:B:577:ARG:HG3	2.07	0.70
1:B:319:ARG:HG2	1:B:592:PHE:HB3	1.72	0.70
1:A:16:VAL:HG21	1:A:258:TRP:HH2	1.58	0.69
1:A:231:ILE:HG22	1:A:233:ILE:HG13	1.73	0.69
1:B:273:ARG:NH1	1:B:290:ASP:OD2	2.24	0.69
1:B:805:ILE:HB	1:B:1054:GLN:NE2	2.08	0.69
1:B:569:ILE:N	1:B:569:ILE:HD12	2.07	0.69
1:C:117:LEU:CD1	1:C:119:ILE:HG13	2.22	0.69
1:A:1054:GLN:HA	1:A:1054:GLN:NE2	2.08	0.69
1:B:456:PHE:HB3	1:B:473:TYR:CD1	2.27	0.69
1:C:392:PHE:HD1	1:C:517:LEU:HD11	1.56	0.68
1:C:176:LEU:HD12	1:C:176:LEU:O	1.93	0.68
1:B:452:LEU:HD21	1:B:492:LEU:HB3	1.74	0.68
1:C:215:ASP:O	1:C:216:LEU:C	2.36	0.68
1:A:319:ARG:HG2	1:A:592:PHE:HB2	1.75	0.68
1:B:112:SER:N	1:B:133:PHE:O	2.27	0.68
1:A:1084:ASP:C	1:A:1086:LYS:H	2.02	0.67
1:C:392:PHE:CD1	1:C:517:LEU:HD11	2.29	0.67
1:C:674:TYR:CE2	1:C:690:GLN:OE1	2.46	0.67
1:A:738:CYS:HB3	1:A:753:LEU:HD21	1.76	0.67
1:B:569:ILE:HD12	1:B:569:ILE:H	1.59	0.67
1:B:159:VAL:HG23	1:B:160:TYR:HD2	1.60	0.67
1:A:347:PHE:CE2	1:A:509:ARG:NH1	2.63	0.67
1:A:849:LEU:O	1:A:853:GLN:HG3	1.95	0.66
1:B:130:VAL:CG1	1:B:168:PHE:HB3	2.25	0.66
1:C:99:ASN:O	1:C:102:ARG:HG3	1.95	0.66
1:C:1054:GLN:HA	1:C:1054:GLN:NE2	2.10	0.66
1:A:422:ASN:HD21	1:A:454:ARG:H	1.43	0.66
1:B:452:LEU:HD21	1:B:492:LEU:CG	2.26	0.66
1:C:103:GLY:HA2	1:C:119:ILE:O	1.96	0.65
1:C:119:ILE:HG12	1:C:128:ILE:HG23	1.77	0.65
1:A:14:GLN:HE22	1:A:156:GLU:HB2	1.62	0.65
1:B:389:ASP:CG	1:B:528:LYS:HD2	2.21	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:492:LEU:HD13	1:C:493:ARG:HD3	1.78	0.65
1:A:88:ASP:O	1:A:270:LEU:HB2	1.98	0.64
1:A:125:ASN:OD1	2:D:1:NAG:C1	2.46	0.64
1:A:347:PHE:CD2	1:A:509:ARG:CZ	2.81	0.64
1:B:456:PHE:CB	1:B:491:PRO:HB3	2.28	0.64
1:A:617:CYS:O	1:A:621:PRO:HD3	1.97	0.63
1:B:456:PHE:N	1:B:491:PRO:HB3	2.14	0.63
1:B:179:LEU:HD23	1:B:188:ASN:HD22	1.63	0.63
1:A:922:LEU:O	1:A:926:GLN:HG3	1.97	0.63
1:C:805:ILE:HB	1:C:1054:GLN:OE1	1.98	0.63
1:B:454:ARG:HG2	1:B:491:PRO:O	1.99	0.63
1:A:1104:VAL:HG23	1:A:1115:ILE:HG12	1.80	0.62
1:B:577:ARG:HG3	1:B:577:ARG:HH11	1.63	0.62
1:C:246:ARG:HG3	1:C:246:ARG:O	1.99	0.62
1:C:998:THR:O	1:C:1002:GLN:HG2	1.99	0.62
1:C:231:ILE:HG23	1:C:231:ILE:O	2.00	0.62
1:C:142:ASP:HB2	1:C:246:ARG:HG2	1.82	0.62
1:C:639:GLY:O	1:C:642:VAL:HG23	1.98	0.62
1:A:1088:HIS:CE1	1:A:1122:VAL:HG22	2.35	0.62
1:B:377:PHE:HE2	1:B:384:PRO:HB3	1.63	0.62
1:B:678:THR:HG22	1:B:678:THR:O	2.00	0.62
1:B:1104:VAL:HG23	1:B:1115:ILE:HG12	1.80	0.62
1:B:441:LEU:HD22	5:B:1403:NAG:H83	1.81	0.61
1:B:452:LEU:CD1	1:B:493:ARG:H	2.14	0.61
1:A:1081:ILE:CD1	1:A:1133:VAL:CG2	2.79	0.61
1:C:677:GLN:N	1:C:677:GLN:OE1	2.32	0.61
1:B:16:VAL:HG22	1:B:16:VAL:O	2.00	0.61
1:B:448:ASN:HB3	1:B:497:PHE:HB2	1.83	0.61
1:C:236:THR:O	1:C:236:THR:HG22	2.01	0.61
1:B:102:ARG:HE	1:B:243:ALA:HB2	1.65	0.60
1:C:357:ARG:NH1	1:C:394:ASN:HA	2.16	0.60
1:B:804:GLN:OE1	1:B:935:GLN:CD	2.44	0.60
1:B:456:PHE:O	1:B:491:PRO:CG	2.47	0.60
1:B:70:VAL:HG12	1:B:70:VAL:O	2.02	0.60
1:A:84:LEU:O	1:A:237:ARG:HG2	2.02	0.60
1:B:58:PHE:HE1	1:B:275:PHE:CE2	2.20	0.60
1:B:677:GLN:OE1	1:B:693:ILE:HG21	2.00	0.60
1:B:190:ARG:HD3	1:B:192:PHE:HE1	1.66	0.59
1:C:332:ILE:HG22	1:C:332:ILE:O	2.02	0.59
1:B:841:LEU:N	1:B:841:LEU:HD23	2.17	0.59
1:B:102:ARG:HD2	1:B:141:LEU:HD12	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:ARG:NH1	1:B:290:ASP:OD1	2.35	0.59
1:A:89:GLY:HA3	1:A:270:LEU:HD12	1.84	0.59
1:B:102:ARG:HE	1:B:243:ALA:CB	2.16	0.59
1:A:16:VAL:HG21	1:A:258:TRP:CH2	2.38	0.58
1:A:752:LEU:HD21	1:A:990:GLU:HG2	1.85	0.58
1:C:577:ARG:HE	1:C:582:LEU:HB3	1.68	0.58
1:C:457:ARG:HD3	1:C:459:SER:O	2.03	0.58
1:C:678:THR:HG23	1:C:678:THR:O	2.01	0.58
1:B:456:PHE:CA	1:B:491:PRO:HB3	2.33	0.58
1:B:452:LEU:CD2	1:B:492:LEU:CD2	2.63	0.58
1:B:569:ILE:H	1:B:569:ILE:CD1	2.16	0.58
1:C:916:LEU:HD12	1:C:923:ILE:HD13	1.86	0.58
1:A:762:GLN:HE21	1:C:961:THR:CG2	2.17	0.58
1:A:755:GLN:OE1	1:C:971:GLY:HA2	2.04	0.58
1:B:452:LEU:CD1	1:B:493:ARG:N	2.64	0.58
1:A:95:THR:O	1:A:95:THR:HG23	2.04	0.58
1:B:1084:ASP:C	1:B:1086:LYS:H	2.11	0.58
1:A:110:LEU:HD12	1:A:110:LEU:N	2.18	0.58
1:B:130:VAL:HG12	1:B:168:PHE:HB3	1.86	0.58
1:B:452:LEU:CG	1:B:492:LEU:HD22	2.31	0.58
1:A:34:ARG:NH1	1:A:221:SER:OG	2.37	0.57
1:B:452:LEU:HD21	1:B:492:LEU:CB	2.34	0.57
1:B:470:THR:CG2	1:B:492:LEU:HD11	2.34	0.57
1:A:517:LEU:HD12	1:A:518:LEU:HG	1.86	0.57
1:B:179:LEU:HD11	1:B:182:LYS:HA	1.86	0.57
1:C:155:SER:HB3	1:C:246:ARG:HH21	1.69	0.57
1:A:357:ARG:NH1	1:A:359:SER:HB3	2.19	0.57
1:C:393:THR:HG21	1:C:519:HIS:HB2	1.87	0.57
1:A:210:ILE:C	1:A:212:LEU:H	2.11	0.57
1:B:1031:GLU:HB2	1:B:1042:PHE:HE2	1.68	0.57
1:C:804:GLN:OE1	2:d:1:NAG:H62	2.05	0.57
1:B:273:ARG:HG3	1:B:273:ARG:HH11	1.68	0.57
1:C:517:LEU:HD12	1:C:517:LEU:N	2.20	0.56
1:B:128:ILE:HD11	1:B:175:PHE:CE1	2.40	0.56
1:B:475:ALA:HB3	1:B:489:TYR:HE1	1.70	0.56
1:C:139:PRO:O	1:C:241:LEU:HB2	2.05	0.56
1:A:83:VAL:HG13	1:A:237:ARG:HD3	1.88	0.56
1:C:126:VAL:HG12	1:C:173:GLN:O	2.05	0.56
1:A:636:TYR:N	1:A:636:TYR:CD1	2.73	0.56
1:B:1093:GLY:O	1:B:1107:ARG:NE	2.24	0.56
1:C:347:PHE:CD1	1:C:509:ARG:HD3	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:ILE:HD11	1:A:235:ILE:HD11	1.87	0.56
1:B:273:ARG:NH1	1:B:290:ASP:CG	2.63	0.56
1:B:377:PHE:CE2	1:B:384:PRO:HB3	2.41	0.56
1:B:489:TYR:N	1:B:489:TYR:CD1	2.73	0.55
1:A:453:TYR:CE2	1:A:455:LEU:HD21	2.40	0.55
1:A:462:LYS:HG2	1:A:465:GLU:HB2	1.87	0.55
1:B:101:ILE:HD11	1:B:240:THR:OG1	2.07	0.55
1:B:319:ARG:HG2	1:B:592:PHE:CB	2.36	0.55
1:B:389:ASP:OD1	1:B:528:LYS:HD2	2.06	0.55
1:C:115:GLN:HB3	1:C:233:ILE:HD13	1.87	0.55
1:C:559:PHE:HZ	1:C:575:ALA:HB3	1.71	0.55
1:B:841:LEU:HD23	1:B:841:LEU:H	1.71	0.55
1:B:139:PRO:HB2	1:B:159:VAL:HG12	1.88	0.55
1:C:18:LEU:HG	1:C:69:HIS:HB2	1.89	0.55
1:C:18:LEU:HG	1:C:18:LEU:O	2.06	0.55
1:B:159:VAL:HG23	1:B:160:TYR:CD2	2.42	0.55
1:B:577:ARG:HH11	1:B:577:ARG:CG	2.18	0.55
1:C:922:LEU:O	1:C:926:GLN:HG3	2.06	0.55
1:A:391:CYS:HA	1:A:525:CYS:HA	1.89	0.54
1:B:246:ARG:HA	1:B:258:TRP:HA	1.90	0.54
1:A:102:ARG:HD2	1:A:141:LEU:HD11	1.88	0.54
1:A:563:GLN:OE1	1:B:43:PHE:HB2	2.07	0.54
1:A:1081:ILE:CD1	1:A:1133:VAL:HG23	2.37	0.54
1:B:96:GLU:OE2	1:B:100:ILE:HB	2.06	0.54
1:B:456:PHE:H	1:B:491:PRO:CB	2.20	0.54
1:C:107:GLY:HA2	1:C:235:ILE:CD1	2.37	0.54
1:C:235:ILE:HG22	1:C:235:ILE:O	2.07	0.54
1:C:231:ILE:HG12	1:C:233:ILE:HG12	1.89	0.54
1:C:639:GLY:O	1:C:642:VAL:CG2	2.55	0.54
1:A:617:CYS:O	1:A:621:PRO:CD	2.54	0.54
1:B:449:TYR:CA	1:B:494:SER:OG	2.56	0.54
1:B:497:PHE:CZ	1:B:507:PRO:HB3	2.43	0.54
1:C:125:ASN:HA	1:C:174:PRO:HB3	1.89	0.54
1:A:819:GLU:HG3	1:A:1054:GLN:NE2	2.23	0.54
1:B:1152:LEU:HD13	1:C:1148:PHE:HE1	1.72	0.54
1:C:498:ARG:HD2	1:C:499:PRO:HD2	1.88	0.54
1:C:117:LEU:HD11	1:C:119:ILE:HG13	1.90	0.54
1:C:120:VAL:HG12	1:C:129:LYS:HZ1	1.73	0.54
1:A:403:ARG:HD2	1:A:505:HIS:NE2	2.23	0.54
1:B:338:PHE:O	1:B:339:ASP:C	2.50	0.54
1:A:395:VAL:HA	1:A:515:PHE:HB3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:444:LYS:HG3	1:C:447:GLY:H	1.73	0.53
1:A:185:ASN:HB3	1:A:212:LEU:HG	1.89	0.53
1:B:497:PHE:CE2	1:B:507:PRO:HB3	2.44	0.53
1:C:1043:CYS:CB	1:C:1048:HIS:CD2	2.92	0.53
1:A:332:ILE:HG22	1:A:362:VAL:HG23	1.89	0.53
2:D:1:NAG:H61	2:D:2:NAG:N2	2.23	0.53
1:B:841:LEU:H	1:B:841:LEU:CD2	2.22	0.53
1:C:1142:GLN:HB2	1:C:1143:PRO:HD3	1.90	0.53
1:C:357:ARG:HH11	1:C:394:ASN:HA	1.73	0.53
1:B:452:LEU:HD21	1:B:492:LEU:HD13	1.91	0.52
1:A:644:GLN:HG2	3:H:1:NAG:H82	1.92	0.52
1:C:836:GLN:C	1:C:838:GLY:H	2.17	0.52
1:A:1081:ILE:HD13	1:A:1133:VAL:HG22	1.92	0.52
1:B:402:ILE:HG22	1:B:403:ARG:O	2.09	0.52
1:A:922:LEU:HD11	1:A:926:GLN:HE21	1.75	0.52
1:B:441:LEU:CD2	5:B:1403:NAG:H83	2.40	0.52
1:A:743:CYS:HB2	1:A:746:SER:HB3	1.92	0.52
1:B:493:ARG:NH1	1:C:378:LYS:NZ	2.58	0.52
1:C:178:ASP:OD2	1:C:182:LYS:HE2	2.10	0.52
1:C:1002:GLN:NE2	1:C:1002:GLN:HA	2.25	0.52
1:B:1093:GLY:HA2	1:B:1107:ARG:HG3	1.91	0.52
1:B:63:THR:CG2	1:B:65:PHE:CE2	2.93	0.51
1:B:454:ARG:HH22	1:B:491:PRO:CG	2.09	0.51
1:C:1043:CYS:HB2	1:C:1048:HIS:CD2	2.45	0.51
1:A:21:ARG:HD2	1:A:21:ARG:O	2.09	0.51
1:B:171:VAL:HG21	2:N:2:NAG:H81	1.92	0.51
1:C:470:THR:HA	1:C:490:PHE:CZ	2.44	0.51
1:C:133:PHE:HB2	1:C:135:PHE:CE2	2.45	0.51
1:A:119:ILE:HD12	1:A:128:ILE:HD12	1.93	0.51
1:C:117:LEU:HD13	1:C:118:LEU:N	2.26	0.51
1:C:1002:GLN:HA	1:C:1002:GLN:HE21	1.76	0.51
1:C:201:PHE:CD1	1:C:201:PHE:C	2.88	0.51
1:A:125:ASN:ND2	2:D:1:NAG:H5	2.26	0.51
1:A:357:ARG:NH1	1:A:359:SER:CB	2.74	0.51
1:C:334:ASN:ND2	1:C:361:CYS:HA	2.25	0.51
1:B:493:ARG:NH1	1:C:378:LYS:HD3	2.04	0.50
1:A:233:ILE:C	1:A:233:ILE:HD12	2.36	0.50
1:B:329:PHE:CE2	1:B:528:LYS:HB2	2.46	0.50
1:C:21:ARG:NH1	1:C:239:GLN:HE22	2.08	0.50
1:A:215:ASP:O	1:A:216:LEU:C	2.55	0.50
1:C:438:SER:OG	1:C:509:ARG:HD2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:GLU:H	1:A:190:ARG:HH22	1.60	0.50
1:B:677:GLN:OE1	1:B:693:ILE:CG2	2.59	0.50
1:B:438:SER:HB2	1:B:509:ARG:HG3	1.93	0.50
1:B:565:PHE:HD1	1:B:576:VAL:HG12	1.76	0.50
1:B:569:ILE:N	1:B:569:ILE:CD1	2.73	0.50
1:B:708:SER:HB3	1:B:711:SER:HB3	1.93	0.50
1:B:445:VAL:O	1:B:498:ARG:NH1	2.46	0.49
1:C:18:LEU:HD21	1:C:69:HIS:HD2	1.77	0.49
1:B:63:THR:HG21	1:B:65:PHE:CE2	2.48	0.49
1:B:805:ILE:HB	1:B:1054:GLN:HE21	1.77	0.49
1:C:115:GLN:HA	1:C:132:GLU:HG2	1.93	0.49
1:C:117:LEU:HD11	1:C:119:ILE:CG1	2.43	0.49
1:A:109:THR:HG22	1:A:114:THR:HB	1.95	0.49
1:A:210:ILE:HG22	1:A:212:LEU:H	1.76	0.49
1:B:456:PHE:HB2	1:B:491:PRO:CB	2.36	0.49
1:B:176:LEU:HD23	1:B:192:PHE:CZ	2.48	0.49
1:C:133:PHE:HB2	1:C:135:PHE:CZ	2.48	0.49
1:A:95:THR:OG1	1:A:187:LYS:HB3	2.13	0.49
1:A:977:LEU:HD11	1:A:993:ILE:HG23	1.95	0.49
1:B:126:VAL:HG13	1:B:174:PRO:HA	1.95	0.49
1:B:498:ARG:HB3	1:B:501:TYR:HB3	1.95	0.49
1:A:58:PHE:HE1	1:A:275:PHE:HE2	1.61	0.48
1:B:426:PRO:HG2	1:B:429:PHE:HB3	1.94	0.48
1:A:58:PHE:CD1	1:A:290:ASP:HB2	2.48	0.48
1:A:125:ASN:HD21	2:D:1:NAG:H3	1.78	0.48
1:A:109:THR:HG1	1:A:111:ASP:CG	2.21	0.48
1:B:1047:TYR:HB2	1:B:1067:TYR:HB3	1.94	0.48
1:B:128:ILE:HD11	1:B:175:PHE:HE1	1.79	0.48
1:C:853:GLN:NE2	1:C:960:ASN:OD1	2.46	0.48
1:B:18:LEU:HD11	1:B:258:TRP:CZ3	2.49	0.48
1:B:676:THR:HG22	1:B:676:THR:O	2.13	0.48
1:A:210:ILE:C	1:A:212:LEU:N	2.72	0.48
1:B:474:GLN:HB2	1:B:488:CYS:SG	2.54	0.48
1:C:708:SER:HB3	1:C:711:SER:HB3	1.95	0.48
1:A:909:ILE:HD13	1:A:1049:LEU:HD21	1.95	0.48
1:A:1081:ILE:HD12	1:A:1133:VAL:HG23	1.96	0.48
1:C:334:ASN:CG	1:C:361:CYS:HA	2.39	0.48
1:A:109:THR:OG1	1:A:111:ASP:OD1	2.31	0.48
1:A:377:PHE:CE2	1:A:384:PRO:HB3	2.42	0.48
1:A:569:ILE:HD12	1:A:569:ILE:H	1.78	0.48
1:B:102:ARG:NE	1:B:243:ALA:HB2	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:475:ALA:CB	1:B:489:TYR:HE1	2.26	0.48
1:C:819:GLU:HG3	1:C:1054:GLN:NE2	2.29	0.48
1:A:347:PHE:CD2	1:A:509:ARG:NH1	2.82	0.47
1:A:897:PRO:HA	1:C:707:TYR:HE2	1.79	0.47
1:C:347:PHE:CE1	1:C:509:ARG:HD3	2.49	0.47
1:A:174:PRO:HG2	1:A:177:MET:SD	2.53	0.47
1:B:273:ARG:NH1	1:B:273:ARG:HG3	2.27	0.47
1:B:1152:LEU:HD13	1:C:1148:PHE:CE1	2.49	0.47
1:C:559:PHE:H	1:C:584:ILE:CD1	2.27	0.47
1:B:493:ARG:HH12	1:C:378:LYS:HD2	1.73	0.47
1:C:1043:CYS:HB3	1:C:1048:HIS:CD2	2.48	0.47
1:A:744:GLY:O	1:A:745:ASP:C	2.58	0.47
1:A:971:GLY:HA3	1:A:995:ARG:NH1	2.26	0.47
1:B:377:PHE:HE2	1:B:384:PRO:CB	2.26	0.47
1:C:456:PHE:HB3	1:C:473:TYR:CG	2.49	0.47
1:B:19:ILE:HD11	1:B:244:LEU:HD11	1.95	0.47
1:B:641:ASN:OD1	1:B:652:GLY:N	2.48	0.47
1:C:347:PHE:CD1	1:C:509:ARG:HB3	2.50	0.47
1:A:347:PHE:CZ	1:A:509:ARG:HG2	2.49	0.47
1:A:1081:ILE:HD13	1:A:1133:VAL:CG2	2.44	0.47
1:A:96:GLU:H	1:A:190:ARG:NH2	2.13	0.47
1:A:490:PHE:HB2	1:A:491:PRO:CD	2.40	0.47
1:B:279:TYR:CE2	1:B:285:ILE:HG13	2.50	0.47
1:B:707:TYR:HE2	1:C:897:PRO:HA	1.79	0.47
1:C:69:HIS:O	1:C:69:HIS:CG	2.67	0.47
1:C:393:THR:HG23	1:C:521:PRO:HD2	1.95	0.47
1:B:470:THR:HG21	1:B:492:LEU:HD11	1.96	0.47
1:B:839:ASP:N	1:B:839:ASP:OD1	2.48	0.47
1:A:21:ARG:CZ	1:A:65:PHE:CE1	2.98	0.47
1:B:442:ASP:O	1:B:448:ASN:ND2	2.48	0.47
1:C:403:ARG:NH2	1:C:495:TYR:HB3	2.30	0.47
1:A:490:PHE:O	1:A:491:PRO:C	2.56	0.46
1:B:452:LEU:HD21	1:B:492:LEU:CD1	2.46	0.46
1:B:475:ALA:HB3	1:B:489:TYR:CE1	2.50	0.46
1:C:99:ASN:O	1:C:102:ARG:CG	2.61	0.46
1:A:42:VAL:HG11	1:C:567:ARG:HD2	1.98	0.46
1:B:19:ILE:HG21	1:B:67:ALA:HB1	1.97	0.46
1:C:231:ILE:CG1	1:C:233:ILE:HG12	2.46	0.46
1:A:84:LEU:O	1:A:237:ARG:CG	2.63	0.46
1:A:365:TYR:HB2	1:A:368:LEU:HD23	1.97	0.46
1:B:22:THR:O	1:B:66:HIS:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:CYS:HB2	1:B:133:PHE:CE2	2.50	0.46
1:B:438:SER:CB	1:B:509:ARG:HG3	2.46	0.46
1:C:120:VAL:O	1:C:120:VAL:HG13	2.15	0.46
1:B:58:PHE:HE1	1:B:275:PHE:HE2	1.59	0.46
1:B:121:ASN:ND2	1:B:174:PRO:HB3	2.27	0.46
1:C:107:GLY:HA2	1:C:235:ILE:HD12	1.97	0.46
1:C:1138:TYR:OH	1:C:1143:PRO:HG2	2.16	0.46
1:A:110:LEU:N	1:A:110:LEU:CD1	2.78	0.46
1:A:371:PHE:HD2	5:A:1403:NAG:H82	1.80	0.46
1:C:456:PHE:HB3	1:C:473:TYR:CD2	2.50	0.46
1:B:338:PHE:O	1:B:341:VAL:HG12	2.16	0.46
1:A:468:ILE:HD13	1:B:115:GLN:HE21	1.81	0.46
1:A:1001:LEU:O	1:A:1005:GLN:HG3	2.15	0.46
1:C:1091:ARG:HD2	1:C:1118:ASP:O	2.15	0.46
1:B:744:GLY:O	1:B:745:ASP:C	2.58	0.46
1:A:130:VAL:O	1:A:130:VAL:HG13	2.16	0.45
1:A:1088:HIS:ND1	1:A:1122:VAL:HG22	2.31	0.45
1:B:319:ARG:HG3	1:B:592:PHE:HB2	1.98	0.45
1:B:320:VAL:HG22	1:B:629:LEU:HD11	1.98	0.45
1:B:830:ASP:C	1:B:832:GLY:H	2.23	0.45
1:A:453:TYR:HE2	1:A:455:LEU:HD21	1.81	0.45
1:A:636:TYR:HD1	1:A:636:TYR:H	1.63	0.45
1:B:102:ARG:CD	1:B:141:LEU:HD12	2.46	0.45
1:B:140:PHE:CD2	1:B:244:LEU:HD13	2.50	0.45
1:B:490:PHE:CG	1:B:490:PHE:O	2.70	0.45
1:A:108:THR:OG1	1:A:234:ASN:O	2.29	0.45
1:A:577:ARG:CZ	1:A:584:ILE:HD11	2.47	0.45
1:A:1142:GLN:HB3	1:A:1143:PRO:HD3	1.99	0.45
1:B:276:LEU:HB3	1:B:289:VAL:CG1	2.46	0.45
1:B:473:TYR:CB	1:B:491:PRO:HD3	2.47	0.45
1:B:905:ARG:HD3	1:B:1049:LEU:O	2.17	0.45
1:C:1081:ILE:HD13	1:C:1133:VAL:HG23	1.97	0.45
1:A:319:ARG:HG2	1:A:592:PHE:CB	2.45	0.45
1:B:1081:ILE:HG12	1:B:1095:PHE:CE2	2.52	0.45
1:A:421:TYR:CD1	1:A:457:ARG:HB2	2.52	0.45
1:B:273:ARG:HD3	1:B:273:ARG:HA	1.71	0.45
1:A:128:ILE:HG21	1:A:229:LEU:HD13	1.98	0.45
1:A:347:PHE:HE2	1:A:509:ARG:NH1	2.15	0.45
1:A:505:HIS:CD2	1:A:505:HIS:O	2.70	0.45
1:A:762:GLN:NE2	1:C:961:THR:CG2	2.80	0.45
1:A:912:THR:OG1	1:A:914:ASN:ND2	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:473:TYR:HB3	1:B:491:PRO:HD3	2.00	0.45
1:B:853:GLN:NE2	1:B:960:ASN:OD1	2.50	0.45
2:G:2:NAG:H4	2:G:3:MAN:H2	1.23	0.45
1:A:567:ARG:NH2	1:A:571:ASP:OD1	2.43	0.44
1:B:473:TYR:HB3	1:B:489:TYR:O	2.17	0.44
1:B:503:VAL:C	1:B:505:HIS:H	2.24	0.44
1:B:21:ARG:HG3	1:B:66:HIS:O	2.17	0.44
1:C:836:GLN:C	1:C:838:GLY:N	2.76	0.44
2:D:1:NAG:H61	2:D:2:NAG:HN2	1.81	0.44
1:B:496:GLY:O	1:B:501:TYR:CE2	2.70	0.44
1:C:922:LEU:HD11	1:C:926:GLN:HE21	1.82	0.44
1:A:33:THR:CG2	1:A:58:PHE:CE2	2.85	0.44
1:B:490:PHE:O	1:B:490:PHE:CD1	2.70	0.44
1:B:677:GLN:CD	1:B:693:ILE:CG2	2.91	0.44
1:C:131:CYS:HB3	1:C:133:PHE:CE2	2.52	0.44
1:A:21:ARG:CZ	1:A:65:PHE:HE1	2.29	0.44
1:A:505:HIS:O	1:A:505:HIS:CG	2.70	0.44
1:B:279:TYR:CE2	1:B:285:ILE:CG1	3.00	0.44
1:C:201:PHE:O	1:C:228:ASP:HA	2.18	0.44
1:B:421:TYR:CE1	1:B:457:ARG:HB3	2.53	0.44
1:A:834:ILE:HD12	1:A:834:ILE:C	2.42	0.44
1:B:830:ASP:C	1:B:832:GLY:N	2.76	0.44
1:B:1048:HIS:HA	1:B:1066:THR:HG22	1.98	0.44
1:C:804:GLN:HA	1:C:804:GLN:NE2	2.33	0.44
1:B:276:LEU:HB3	1:B:289:VAL:HG12	1.99	0.43
1:B:493:ARG:CZ	1:C:378:LYS:HD2	2.47	0.43
1:A:96:GLU:HA	1:A:96:GLU:OE2	2.18	0.43
1:A:338:PHE:HZ	1:A:367:VAL:HG23	1.83	0.43
1:A:636:TYR:N	1:A:636:TYR:HD1	2.15	0.43
1:A:403:ARG:NE	1:A:505:HIS:CE1	2.86	0.43
1:B:739:THR:O	1:B:743:CYS:O	2.36	0.43
1:C:490:PHE:HB3	1:C:491:PRO:HD3	2.00	0.43
1:B:183:GLN:HG2	1:B:184:GLY:N	2.33	0.43
1:B:456:PHE:HB3	1:B:473:TYR:HD1	1.78	0.43
1:A:347:PHE:CE2	1:A:509:ARG:HG2	2.54	0.43
1:C:517:LEU:CD1	1:C:517:LEU:N	2.81	0.43
1:A:346:ARG:HA	1:A:509:ARG:NH2	2.34	0.43
1:B:176:LEU:HD13	1:B:190:ARG:NH1	2.34	0.43
1:B:569:ILE:C	1:B:571:ASP:H	2.26	0.43
1:C:141:LEU:HB2	1:C:241:LEU:HD13	2.00	0.43
1:C:244:LEU:HD13	1:C:244:LEU:HA	1.75	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:805:ILE:HB	1:A:1054:GLN:OE1	2.18	0.43
1:A:1054:GLN:HE21	1:A:1054:GLN:CA	2.17	0.43
1:C:140:PHE:CD2	1:C:140:PHE:C	2.97	0.43
1:C:14:GLN:HE22	1:C:158:ARG:HA	1.83	0.43
1:C:1153:ASP:OD2	1:C:1157:LYS:HE2	2.19	0.43
1:A:431:GLY:HA2	1:A:515:PHE:CE2	2.53	0.43
1:B:490:PHE:CE2	1:B:492:LEU:HD12	2.53	0.43
1:C:347:PHE:CE1	1:C:509:ARG:HB3	2.54	0.43
1:A:355:ARG:NH1	1:A:398:ASP:OD1	2.52	0.42
1:B:377:PHE:HE2	1:B:384:PRO:CG	2.31	0.42
1:B:493:ARG:NH1	1:C:378:LYS:HD2	2.29	0.42
1:B:493:ARG:NH1	1:C:378:LYS:HZ3	2.15	0.42
1:B:752:LEU:HD11	1:B:990:GLU:HG3	2.01	0.42
1:C:1084:ASP:C	1:C:1086:LYS:H	2.27	0.42
1:C:1141:LEU:HD13	1:C:1141:LEU:C	2.44	0.42
1:A:18:LEU:HG	1:A:69:HIS:CE1	2.54	0.42
1:B:470:THR:HG22	1:B:492:LEU:HD11	2.00	0.42
1:B:708:SER:HB3	1:B:711:SER:CB	2.49	0.42
1:A:414:GLN:OE1	1:A:414:GLN:HA	2.19	0.42
1:C:121:ASN:HD22	1:C:176:LEU:HG	1.84	0.42
1:A:18:LEU:HD11	1:A:70:VAL:O	2.19	0.42
1:B:16:VAL:HG22	1:B:18:LEU:HG	2.00	0.42
1:C:599:THR:HB	1:C:608:VAL:HG12	2.01	0.42
1:C:868:GLU:O	1:C:872:GLN:HG3	2.20	0.42
1:A:392:PHE:HB2	1:A:395:VAL:HG23	2.02	0.42
1:B:205:SER:OG	1:B:226:LEU:HD21	2.20	0.42
1:C:127:VAL:HG21	2:X:1:NAG:O6	2.19	0.42
1:A:144:TYR:CD1	1:A:155:SER:HB2	2.54	0.42
1:B:805:ILE:CB	1:B:1054:GLN:NE2	2.82	0.42
1:C:22:THR:HG23	1:C:66:HIS:HB2	2.02	0.42
1:C:417:ASN:CG	1:C:455:LEU:HD22	2.45	0.42
1:C:495:TYR:HB2	1:C:497:PHE:CZ	2.55	0.42
1:C:1139:ASP:HB2	1:C:1142:GLN:HG2	2.00	0.42
1:B:246:ARG:HA	1:B:258:TRP:CD1	2.55	0.42
1:B:179:LEU:HD23	1:B:188:ASN:ND2	2.31	0.42
1:B:319:ARG:CG	1:B:592:PHE:HB2	2.50	0.42
1:C:69:HIS:O	1:C:69:HIS:ND1	2.52	0.42
1:A:209:PRO:O	1:A:210:ILE:HD13	2.20	0.42
1:C:121:ASN:OD1	1:C:121:ASN:C	2.62	0.42
1:A:748:GLU:H	1:A:748:GLU:CD	2.27	0.42
1:B:176:LEU:HD13	1:B:190:ARG:HH12	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:490:PHE:HD1	1:B:490:PHE:H	1.67	0.42
1:C:101:ILE:HD12	1:C:241:LEU:O	2.20	0.41
1:C:357:ARG:HG3	1:C:395:VAL:O	2.20	0.41
1:B:630:THR:HB	1:B:631:PRO:HD3	2.02	0.41
1:B:841:LEU:N	1:B:841:LEU:CD2	2.80	0.41
1:C:320:VAL:HG13	1:C:590:CYS:HB3	2.02	0.41
1:B:406:GLU:C	1:B:408:SER:N	2.78	0.41
1:A:101:ILE:HD13	1:A:242:LEU:HD11	2.03	0.41
1:A:197:ILE:O	1:A:197:ILE:HG23	2.20	0.41
1:A:233:ILE:C	1:A:233:ILE:CD1	2.93	0.41
1:C:229:LEU:HA	1:C:230:PRO:HD2	1.84	0.41
1:C:759:PHE:O	1:C:763:LEU:HD22	2.20	0.41
1:A:41:LYS:CE	1:C:562:PHE:O	2.68	0.41
1:B:21:ARG:HA	1:B:66:HIS:O	2.20	0.41
1:B:319:ARG:CG	1:B:592:PHE:CB	2.99	0.41
1:B:655:TYR:HA	1:B:694:ALA:O	2.19	0.41
1:C:69:HIS:HE1	1:C:261:GLY:N	2.19	0.41
1:C:241:LEU:HD12	1:C:242:LEU:N	2.35	0.41
1:B:244:LEU:HB3	1:B:258:TRP:CD1	2.56	0.41
1:B:849:LEU:O	1:B:852:ALA:HB3	2.20	0.41
1:B:1144:GLU:H	1:B:1144:GLU:HG3	1.67	0.41
1:C:111:ASP:CG	1:C:113:LYS:HD3	2.46	0.41
1:A:176:LEU:HA	1:A:207:HIS:CE1	2.56	0.41
1:A:425:LEU:HD23	1:A:425:LEU:H	1.85	0.41
1:A:980:ILE:HG21	1:A:992:GLN:HB2	2.03	0.41
2:N:1:NAG:H61	2:N:2:NAG:N2	2.36	0.41
1:A:109:THR:CG2	1:A:111:ASP:OD1	2.69	0.41
1:B:176:LEU:HB2	1:B:190:ARG:CZ	2.50	0.41
1:B:338:PHE:CE2	1:B:513:LEU:HD11	2.56	0.41
1:B:454:ARG:HH12	1:B:491:PRO:CG	2.23	0.41
1:C:139:PRO:O	1:C:241:LEU:CB	2.68	0.41
1:C:347:PHE:HD1	1:C:509:ARG:HD3	1.85	0.41
1:A:18:LEU:HD23	1:A:69:HIS:CE1	2.56	0.41
1:A:21:ARG:C	1:A:21:ARG:CD	2.94	0.41
1:A:41:LYS:HE3	1:C:562:PHE:O	2.21	0.41
1:A:470:THR:HG22	1:A:491:PRO:HG2	2.03	0.41
1:B:347:PHE:CD2	1:B:509:ARG:HD2	2.56	0.41
1:C:337:PRO:O	1:C:340:GLU:HG2	2.21	0.41
1:A:139:PRO:O	1:A:241:LEU:CD1	2.70	0.40
1:B:183:GLN:HB3	1:B:186:PHE:CB	2.51	0.40
1:B:1102:TRP:CZ2	1:B:1133:VAL:HG21	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1141:LEU:O	1:C:1144:GLU:HG3	2.21	0.40
1:A:708:SER:HB3	1:A:711:SER:HB3	2.03	0.40
1:A:1126:CYS:HB2	1:A:1132:ILE:HD13	2.03	0.40
1:B:359:SER:HA	1:B:524:VAL:CG2	2.51	0.40
1:C:69:HIS:CE1	1:C:260:ALA:HA	2.56	0.40
1:A:246:ARG:HA	1:A:258:TRP:H	1.86	0.40
1:A:762:GLN:HE21	1:C:961:THR:HG21	1.87	0.40
1:B:124:THR:O	1:B:174:PRO:HD3	2.21	0.40
1:B:490:PHE:CE2	1:B:492:LEU:HB2	2.56	0.40
1:B:1142:GLN:HB2	1:B:1143:PRO:HD3	2.02	0.40
1:C:708:SER:HB3	1:C:711:SER:CB	2.51	0.40
1:C:874:THR:HA	1:C:877:LEU:HD12	2.03	0.40
1:A:21:ARG:NH2	1:A:65:PHE:CE1	2.90	0.40
1:A:33:THR:HA	1:A:58:PHE:CE2	2.57	0.40
1:A:1084:ASP:C	1:A:1086:LYS:N	2.70	0.40
1:B:112:SER:CA	1:B:133:PHE:O	2.70	0.40
1:C:266:TYR:CD1	1:C:266:TYR:N	2.88	0.40
1:C:271:GLN:HE21	1:C:271:GLN:HB2	1.66	0.40
1:C:1101:HIS:ND1	2:f:1:NAG:H5	2.36	0.40
1:A:133:PHE:HB2	1:A:135:PHE:CE2	2.57	0.40
1:B:449:TYR:C	1:B:494:SER:HG	2.28	0.40
1:C:239:GLN:HA	1:C:239:GLN:OE1	2.22	0.40
1:C:335:LEU:HD13	1:C:336:CYS:N	2.37	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	1100/1305 (84%)	1010 (92%)	88 (8%)	2 (0%)	43	76
1	B	1116/1305 (86%)	1045 (94%)	70 (6%)	1 (0%)	48	80

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	C	1111/1305 (85%)	1014 (91%)	96 (9%)	1 (0%)	48 80
All	All	3327/3915 (85%)	3069 (92%)	254 (8%)	4 (0%)	49 80

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	155	SER
1	C	520	ALA
1	A	491	PRO
1	B	491	PRO

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	968/1130 (86%)	951 (98%)	17 (2%)	51 77
1	B	978/1130 (86%)	950 (97%)	28 (3%)	37 70
1	C	974/1130 (86%)	943 (97%)	31 (3%)	34 67
All	All	2920/3390 (86%)	2844 (97%)	76 (3%)	41 72

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	144	TYR
1	A	159	VAL
1	A	212	LEU
1	A	368	LEU
1	A	394	ASN
1	A	402	ILE
1	A	461	LEU
1	A	491	PRO
1	A	492	LEU
1	A	495	TYR
1	A	506	GLN

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Mol	Chain	Res	Type
1	A	509	ARG
1	A	524	VAL
1	A	676	THR
1	A	914	ASN
1	A	916	LEU
1	A	1050	MET
1	B	16	VAL
1	B	117	LEU
1	B	118	LEU
1	B	154	GLU
1	B	176	LEU
1	B	177	MET
1	B	179	LEU
1	B	197	ILE
1	B	229	LEU
1	B	233	ILE
1	B	240	THR
1	B	245	HIS
1	B	273	ARG
1	B	323	THR
1	B	364	ASP
1	B	453	TYR
1	B	461	LEU
1	B	466	ARG
1	B	486	PHE
1	B	489	TYR
1	B	567	ARG
1	B	569	ILE
1	B	581	THR
1	B	629	LEU
1	B	634	ARG
1	B	774	GLN
1	B	776	LYS
1	B	779	GLN
1	C	69	HIS
1	C	70	VAL
1	C	81	ASN
1	C	108	THR
1	C	126	VAL
1	C	130	VAL
1	C	140	PHE
1	C	153	MET

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Mol	Chain	Res	Type
1	C	154	GLU
1	C	186	PHE
1	C	189	LEU
1	C	201	PHE
1	C	238	PHE
1	C	244	LEU
1	C	266	TYR
1	C	271	GLN
1	C	318	PHE
1	C	320	VAL
1	C	356	LYS
1	C	358	ILE
1	C	366	SER
1	C	371	PHE
1	C	375	PHE
1	C	425	LEU
1	C	511	VAL
1	C	569	ILE
1	C	599	THR
1	C	763	LEU
1	C	878	LEU
1	C	914	ASN
1	C	916	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (69) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	81	ASN
1	A	99	ASN
1	A	121	ASN
1	A	173	GLN
1	A	196	ASN
1	A	218	GLN
1	A	317	ASN
1	A	321	GLN
1	A	394	ASN
1	A	422	ASN
1	A	437	ASN
1	A	450	ASN
1	A	474	GLN
1	A	487	ASN

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Mol	Chain	Res	Type
1	A	519	HIS
1	A	606	ASN
1	A	703	ASN
1	A	762	GLN
1	A	913	GLN
1	A	914	ASN
1	A	954	HIS
1	A	1054	GLN
1	A	1083	HIS
1	A	1106	GLN
1	A	1142	GLN
1	A	1159	HIS
1	B	30	ASN
1	B	52	GLN
1	B	69	HIS
1	B	81	ASN
1	B	115	GLN
1	B	121	ASN
1	B	149	ASN
1	B	173	GLN
1	B	188	ASN
1	B	207	HIS
1	B	370	ASN
1	B	394	ASN
1	B	439	ASN
1	B	506	GLN
1	B	536	ASN
1	B	544	ASN
1	B	564	GLN
1	B	613	GLN
1	B	751	ASN
1	B	853	GLN
1	B	1005	GLN
1	B	1071	GLN
1	B	1088	HIS
1	B	1142	GLN
1	C	14	GLN
1	C	69	HIS
1	C	81	ASN
1	C	149	ASN
1	C	218	GLN
1	C	239	GLN

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Mol	Chain	Res	Type
1	C	271	GLN
1	C	314	GLN
1	C	321	GLN
1	C	460	ASN
1	C	606	ASN
1	C	607	GLN
1	C	856	ASN
1	C	872	GLN
1	C	913	GLN
1	C	1002	GLN
1	C	1054	GLN
1	C	1159	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

81 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	0.33	0	17,19,21	0.49	0
2	NAG	D	2	2	14,14,15	0.20	0	17,19,21	0.50	0
2	MAN	D	3	2	11,11,12	0.73	0	15,15,17	0.99	2 (13%)
3	NAG	E	1	3,1	14,14,15	0.33	0	17,19,21	0.45	0
3	NAG	E	2	3	14,14,15	0.20	0	17,19,21	0.44	0
3	NAG	F	1	3,1	14,14,15	0.39	0	17,19,21	0.32	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	F	2	3	14,14,15	0.40	0	17,19,21	0.37	0
2	NAG	G	1	2,1	14,14,15	0.40	0	17,19,21	0.35	0
2	NAG	G	2	2	14,14,15	0.40	0	17,19,21	0.40	0
2	MAN	G	3	2	11,11,12	0.26	0	15,15,17	0.55	0
3	NAG	H	1	3,1	14,14,15	0.40	0	17,19,21	0.50	0
3	NAG	H	2	3	14,14,15	0.40	0	17,19,21	0.37	0
2	NAG	I	1	2,1	14,14,15	0.26	0	17,19,21	0.46	0
2	NAG	I	2	2	14,14,15	0.22	0	17,19,21	0.42	0
2	MAN	I	3	2	11,11,12	0.72	0	15,15,17	1.26	2 (13%)
2	NAG	J	1	2,1	14,14,15	0.32	0	17,19,21	0.45	0
2	NAG	J	2	2	14,14,15	0.22	0	17,19,21	0.42	0
2	MAN	J	3	2	11,11,12	0.67	0	15,15,17	1.04	2 (13%)
4	NAG	K	1	4,1	14,14,15	0.21	0	17,19,21	0.85	1 (5%)
4	NAG	K	2	4	14,14,15	0.25	0	17,19,21	0.42	0
4	FUC	K	3	4	10,10,11	0.99	0	14,14,16	1.02	2 (14%)
2	NAG	L	1	2,1	14,14,15	0.21	0	17,19,21	0.48	0
2	NAG	L	2	2	14,14,15	0.20	0	17,19,21	0.43	0
2	MAN	L	3	2	11,11,12	0.80	0	15,15,17	1.01	2 (13%)
2	NAG	M	1	2,1	14,14,15	0.26	0	17,19,21	0.49	0
2	NAG	M	2	2	14,14,15	0.20	0	17,19,21	0.47	0
2	MAN	M	3	2	11,11,12	0.74	0	15,15,17	1.33	1 (6%)
2	NAG	N	1	2,1	14,14,15	0.42	0	17,19,21	0.50	0
2	NAG	N	2	2	14,14,15	0.26	0	17,19,21	0.54	0
2	MAN	N	3	2	11,11,12	0.65	0	15,15,17	1.25	2 (13%)
3	NAG	O	1	3,1	14,14,15	0.41	0	17,19,21	0.43	0
3	NAG	O	2	3	14,14,15	0.25	0	17,19,21	0.45	0
3	NAG	P	1	3,1	14,14,15	0.39	0	17,19,21	0.42	0
3	NAG	P	2	3	14,14,15	0.41	0	17,19,21	0.50	0
2	NAG	Q	1	2,1	14,14,15	0.24	0	17,19,21	0.40	0
2	NAG	Q	2	2	14,14,15	0.24	0	17,19,21	0.45	0
2	MAN	Q	3	2	11,11,12	0.68	0	15,15,17	1.21	2 (13%)
3	NAG	R	1	3,1	14,14,15	0.39	0	17,19,21	0.50	0
3	NAG	R	2	3	14,14,15	0.39	0	17,19,21	0.37	0
2	NAG	S	1	2,1	14,14,15	0.29	0	17,19,21	0.49	0
2	NAG	S	2	2	14,14,15	0.19	0	17,19,21	0.42	0
2	MAN	S	3	2	11,11,12	0.73	0	15,15,17	1.27	2 (13%)
2	NAG	T	1	2,1	14,14,15	0.22	0	17,19,21	0.48	0
2	NAG	T	2	2	14,14,15	0.18	0	17,19,21	0.48	0
2	MAN	T	3	2	11,11,12	0.68	0	15,15,17	0.95	2 (13%)
4	NAG	U	1	4,1	14,14,15	0.36	0	17,19,21	0.79	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	U	2	4	14,14,15	0.21	0	17,19,21	0.42	0
4	FUC	U	3	4	10,10,11	0.85	0	14,14,16	1.03	2 (14%)
2	NAG	V	1	2,1	14,14,15	0.23	0	17,19,21	0.46	0
2	NAG	V	2	2	14,14,15	0.19	0	17,19,21	0.45	0
2	MAN	V	3	2	11,11,12	0.80	0	15,15,17	1.13	2 (13%)
2	NAG	W	1	2,1	14,14,15	0.25	0	17,19,21	0.49	0
2	NAG	W	2	2	14,14,15	0.24	0	17,19,21	0.51	0
2	MAN	W	3	2	11,11,12	0.67	0	15,15,17	0.97	2 (13%)
2	NAG	X	1	2,1	14,14,15	0.39	0	17,19,21	1.08	1 (5%)
2	NAG	X	2	2	14,14,15	0.22	0	17,19,21	0.52	0
2	MAN	X	3	2	11,11,12	0.73	0	15,15,17	1.24	2 (13%)
3	NAG	Y	1	3,1	14,14,15	0.28	0	17,19,21	0.39	0
3	NAG	Y	2	3	14,14,15	0.24	0	17,19,21	0.45	0
3	NAG	Z	1	3,1	14,14,15	0.28	0	17,19,21	0.52	0
3	NAG	Z	2	3	14,14,15	0.18	0	17,19,21	0.51	0
2	NAG	a	1	2,1	14,14,15	0.25	0	17,19,21	0.42	0
2	NAG	a	2	2	14,14,15	0.19	0	17,19,21	0.46	0
2	MAN	a	3	2	11,11,12	0.68	0	15,15,17	1.06	2 (13%)
3	NAG	b	1	3,1	14,14,15	0.41	0	17,19,21	0.99	2 (11%)
3	NAG	b	2	3	14,14,15	0.40	0	17,19,21	0.39	0
2	NAG	c	1	2,1	14,14,15	0.24	0	17,19,21	0.50	0
2	NAG	c	2	2	14,14,15	0.23	0	17,19,21	0.41	0
2	MAN	c	3	2	11,11,12	0.69	0	15,15,17	1.33	2 (13%)
2	NAG	d	1	2,1	14,14,15	0.30	0	17,19,21	0.41	0
2	NAG	d	2	2	14,14,15	0.19	0	17,19,21	0.47	0
2	MAN	d	3	2	11,11,12	0.72	0	15,15,17	1.08	2 (13%)
4	NAG	e	1	4,1	14,14,15	0.19	0	17,19,21	0.72	1 (5%)
4	NAG	e	2	4	14,14,15	0.25	0	17,19,21	0.40	0
4	FUC	e	3	4	10,10,11	0.92	0	14,14,16	0.92	1 (7%)
2	NAG	f	1	2,1	14,14,15	0.26	0	17,19,21	0.48	0
2	NAG	f	2	2	14,14,15	0.21	0	17,19,21	0.50	0
2	MAN	f	3	2	11,11,12	0.79	0	15,15,17	1.17	2 (13%)
2	NAG	g	1	2,1	14,14,15	0.21	0	17,19,21	0.47	0
2	NAG	g	2	2	14,14,15	0.20	0	17,19,21	0.43	0
2	MAN	g	3	2	11,11,12	0.79	0	15,15,17	1.44	2 (13%)

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	-	2/6/23/26	0/1/1/1
2	MAN	D	3	2	-	1/2/19/22	1/1/1/1
3	NAG	E	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	NAG	F	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	MAN	G	3	2	-	0/2/19/22	1/1/1/1
3	NAG	H	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
2	NAG	I	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	2/6/23/26	0/1/1/1
2	MAN	I	3	2	-	2/2/19/22	1/1/1/1
2	NAG	J	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	MAN	J	3	2	-	2/2/19/22	1/1/1/1
4	NAG	K	1	4,1	-	1/6/23/26	0/1/1/1
4	NAG	K	2	4	-	0/6/23/26	0/1/1/1
4	FUC	K	3	4	-	-	0/1/1/1
2	NAG	L	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	L	2	2	-	0/6/23/26	0/1/1/1
2	MAN	L	3	2	-	1/2/19/22	1/1/1/1
2	NAG	M	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	M	2	2	-	2/6/23/26	0/1/1/1
2	MAN	M	3	2	-	0/2/19/22	1/1/1/1
2	NAG	N	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	N	2	2	-	2/6/23/26	0/1/1/1
2	MAN	N	3	2	-	1/2/19/22	1/1/1/1
3	NAG	O	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	O	2	3	-	2/6/23/26	0/1/1/1
3	NAG	P	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	P	2	3	-	0/6/23/26	0/1/1/1
2	NAG	Q	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	Q	2	2	-	0/6/23/26	0/1/1/1
2	MAN	Q	3	2	-	0/2/19/22	1/1/1/1
3	NAG	R	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	R	2	3	-	0/6/23/26	0/1/1/1
2	NAG	S	1	2,1	-	1/6/23/26	0/1/1/1

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	g	2	2	-	2/6/23/26	0/1/1/1
2	MAN	g	3	2	-	0/2/19/22	1/1/1/1

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	g	3	MAN	C1-O5-C5	4.70	118.49	112.19
2	M	3	MAN	C1-O5-C5	4.35	118.01	112.19
2	c	3	MAN	C1-O5-C5	4.25	117.88	112.19
2	S	3	MAN	C1-O5-C5	3.96	117.49	112.19
2	N	3	MAN	C1-O5-C5	3.95	117.48	112.19
2	I	3	MAN	C1-O5-C5	3.92	117.44	112.19
2	X	1	NAG	O5-C1-C2	-3.85	105.33	111.29
2	X	3	MAN	C1-O5-C5	3.85	117.35	112.19
2	Q	3	MAN	C1-O5-C5	3.68	117.12	112.19
2	f	3	MAN	C1-O5-C5	3.43	116.79	112.19
2	V	3	MAN	C1-O5-C5	3.29	116.60	112.19
4	K	1	NAG	C1-O5-C5	3.14	116.39	112.19
2	d	3	MAN	C1-O5-C5	3.11	116.36	112.19
2	a	3	MAN	C1-O5-C5	3.03	116.25	112.19
2	J	3	MAN	C1-O5-C5	2.97	116.17	112.19
3	b	1	NAG	C2-N2-C7	2.80	126.66	122.90
4	U	1	NAG	C1-O5-C5	2.78	115.91	112.19
2	L	3	MAN	C1-O5-C5	2.76	115.89	112.19
2	W	3	MAN	C1-O5-C5	2.64	115.72	112.19
2	D	3	MAN	C1-O5-C5	2.62	115.70	112.19
4	e	1	NAG	C1-O5-C5	2.54	115.58	112.19
2	T	3	MAN	C1-O5-C5	2.49	115.53	112.19
2	Q	3	MAN	O2-C2-C3	-2.29	105.40	110.15
3	b	1	NAG	C1-C2-N2	2.29	114.05	110.43
2	f	3	MAN	O2-C2-C3	-2.28	105.42	110.15
2	D	3	MAN	O2-C2-C3	-2.25	105.48	110.15
2	S	3	MAN	O2-C2-C3	-2.23	105.53	110.15
2	V	3	MAN	O2-C2-C3	-2.21	105.58	110.15
2	X	3	MAN	O2-C2-C3	-2.20	105.59	110.15
2	J	3	MAN	O2-C2-C3	-2.20	105.60	110.15
2	d	3	MAN	O2-C2-C3	-2.20	105.60	110.15
2	c	3	MAN	O2-C2-C3	-2.19	105.61	110.15
2	a	3	MAN	O2-C2-C3	-2.18	105.63	110.15
4	K	3	FUC	C1-C2-C3	2.18	112.81	109.64
2	T	3	MAN	O2-C2-C3	-2.17	105.65	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	3	MAN	O2-C2-C3	-2.14	105.71	110.15
2	N	3	MAN	O2-C2-C3	-2.14	105.72	110.15
2	L	3	MAN	O2-C2-C3	-2.14	105.73	110.15
2	W	3	MAN	O2-C2-C3	-2.12	105.77	110.15
4	K	3	FUC	O5-C5-C4	2.11	113.35	109.55
4	U	3	FUC	C1-O5-C5	2.06	117.83	112.97
4	U	3	FUC	O5-C5-C4	2.05	113.24	109.55
4	e	3	FUC	O5-C5-C4	2.03	113.20	109.55
2	g	3	MAN	O2-C2-C3	-2.01	105.98	110.15

There are no chirality outliers.

All (62) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	b	1	NAG	C1-C2-N2-C7
3	Z	1	NAG	O5-C5-C6-O6
2	g	1	NAG	O5-C5-C6-O6
2	J	3	MAN	O5-C5-C6-O6
2	f	3	MAN	O5-C5-C6-O6
2	g	1	NAG	C4-C5-C6-O6
3	Z	1	NAG	C4-C5-C6-O6
2	V	3	MAN	O5-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	M	2	NAG	O5-C5-C6-O6
2	T	1	NAG	O5-C5-C6-O6
2	J	1	NAG	O5-C5-C6-O6
2	S	2	NAG	O5-C5-C6-O6
2	a	1	NAG	O5-C5-C6-O6
2	D	1	NAG	O5-C5-C6-O6
4	U	1	NAG	O5-C5-C6-O6
2	T	1	NAG	C4-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	J	1	NAG	C4-C5-C6-O6
2	M	2	NAG	C4-C5-C6-O6
2	a	1	NAG	C4-C5-C6-O6
2	I	2	NAG	O5-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
2	I	3	MAN	C4-C5-C6-O6
2	X	2	NAG	C4-C5-C6-O6
2	N	2	NAG	O5-C5-C6-O6
2	D	3	MAN	O5-C5-C6-O6
2	L	3	MAN	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	O	2	NAG	C4-C5-C6-O6
3	O	2	NAG	O5-C5-C6-O6
2	N	2	NAG	C4-C5-C6-O6
2	J	3	MAN	C4-C5-C6-O6
2	f	3	MAN	C4-C5-C6-O6
3	Z	2	NAG	C4-C5-C6-O6
3	Z	2	NAG	O5-C5-C6-O6
2	X	2	NAG	O5-C5-C6-O6
4	K	1	NAG	O5-C5-C6-O6
2	X	3	MAN	O5-C5-C6-O6
2	d	3	MAN	O5-C5-C6-O6
2	N	3	MAN	O5-C5-C6-O6
2	N	1	NAG	C4-C5-C6-O6
2	d	2	NAG	C4-C5-C6-O6
2	d	2	NAG	O5-C5-C6-O6
3	b	1	NAG	C3-C2-N2-C7
2	T	3	MAN	O5-C5-C6-O6
2	S	2	NAG	C4-C5-C6-O6
4	U	1	NAG	C4-C5-C6-O6
2	S	1	NAG	O5-C5-C6-O6
2	N	1	NAG	O5-C5-C6-O6
2	V	3	MAN	C4-C5-C6-O6
2	g	2	NAG	C4-C5-C6-O6
2	I	3	MAN	O5-C5-C6-O6
2	J	2	NAG	C4-C5-C6-O6
2	I	2	NAG	C4-C5-C6-O6
3	Y	1	NAG	O5-C5-C6-O6
2	V	1	NAG	C4-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
2	c	2	NAG	O5-C5-C6-O6
4	e	1	NAG	O5-C5-C6-O6
2	g	2	NAG	O5-C5-C6-O6
2	V	1	NAG	O5-C5-C6-O6
3	R	1	NAG	C3-C2-N2-C7

All (18) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	3	MAN	C1-C2-C3-C4-C5-O5
2	T	3	MAN	C1-C2-C3-C4-C5-O5
2	D	3	MAN	C1-C2-C3-C4-C5-O5
2	g	3	MAN	C1-C2-C3-C4-C5-O5

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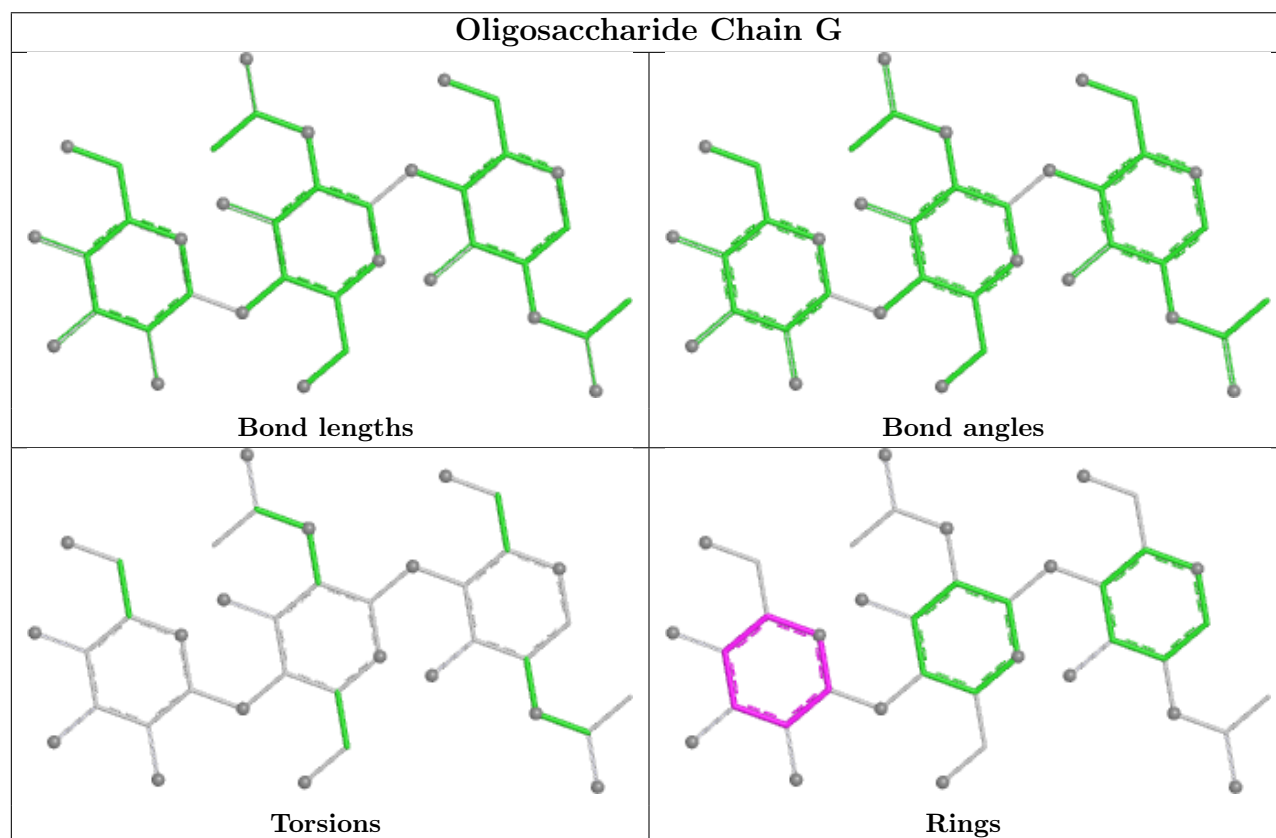
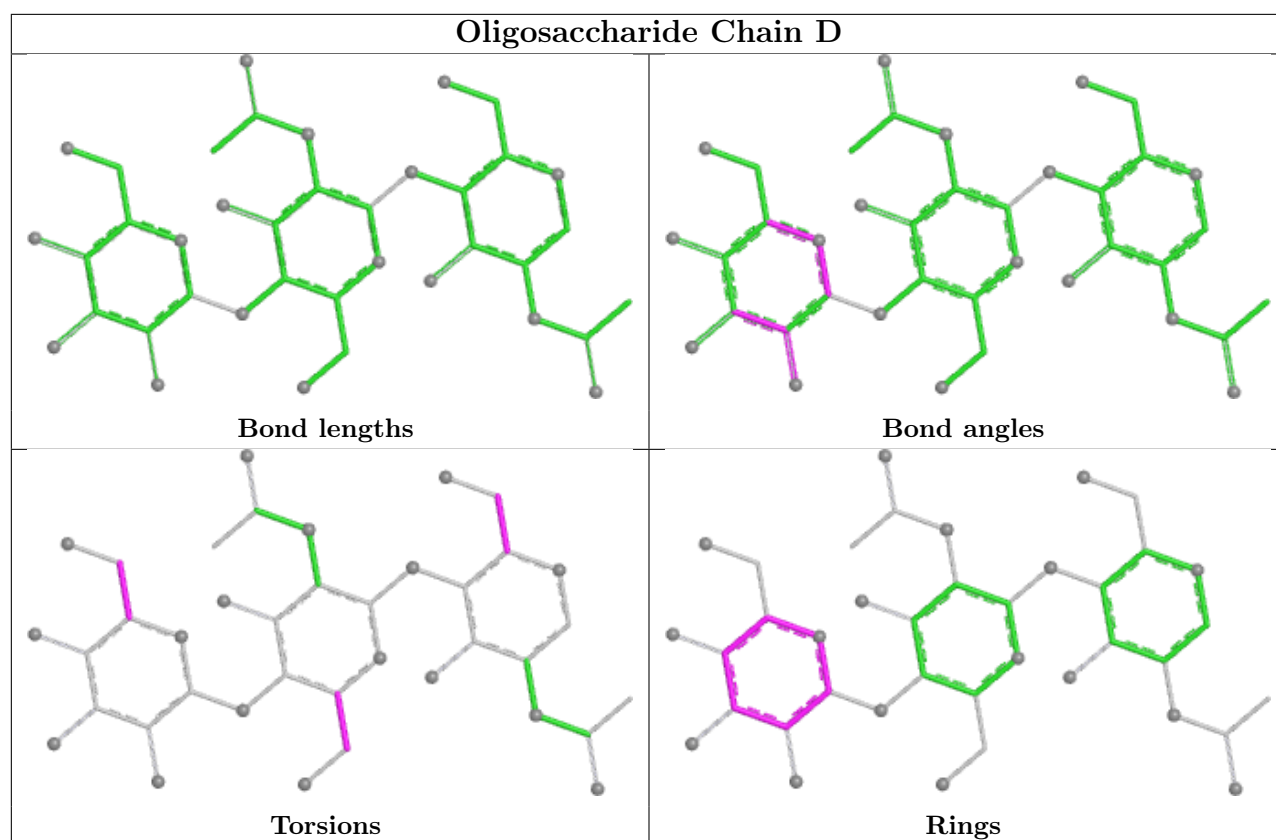
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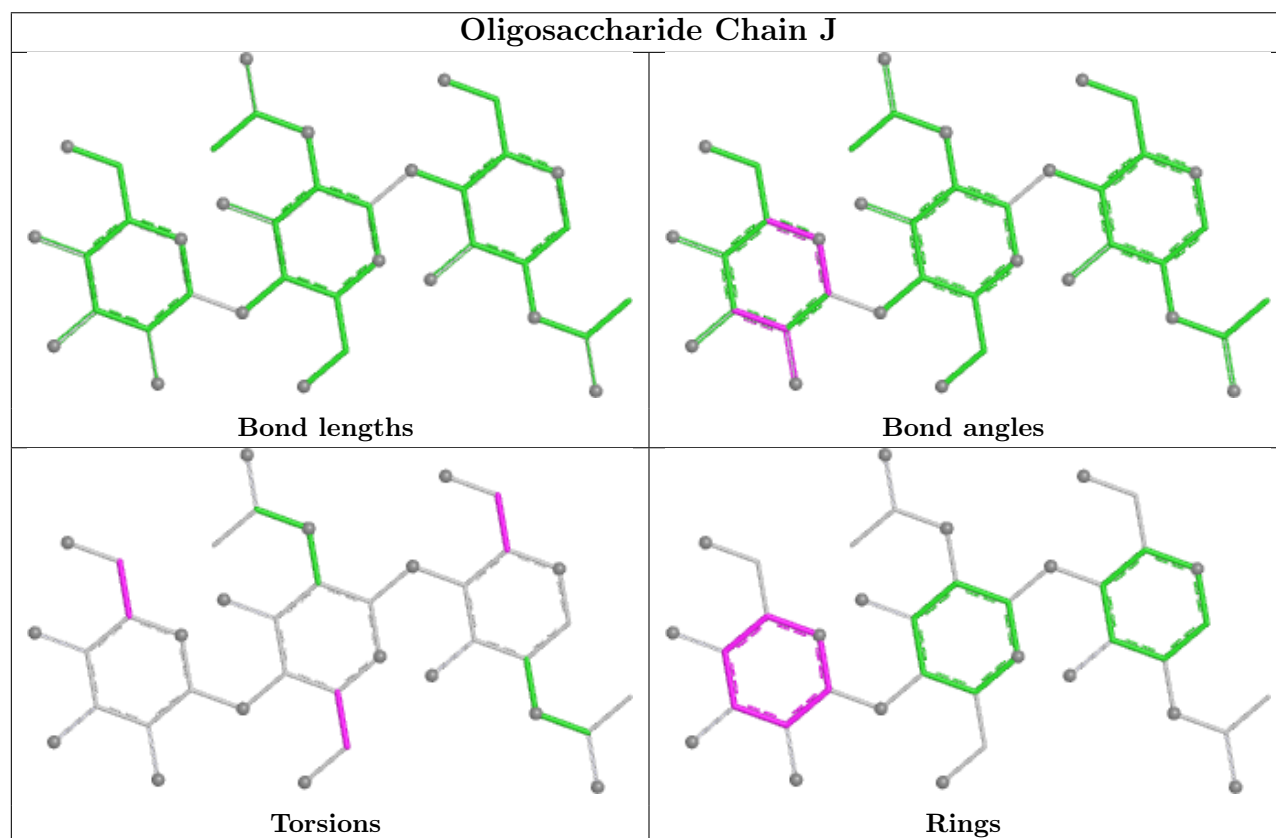
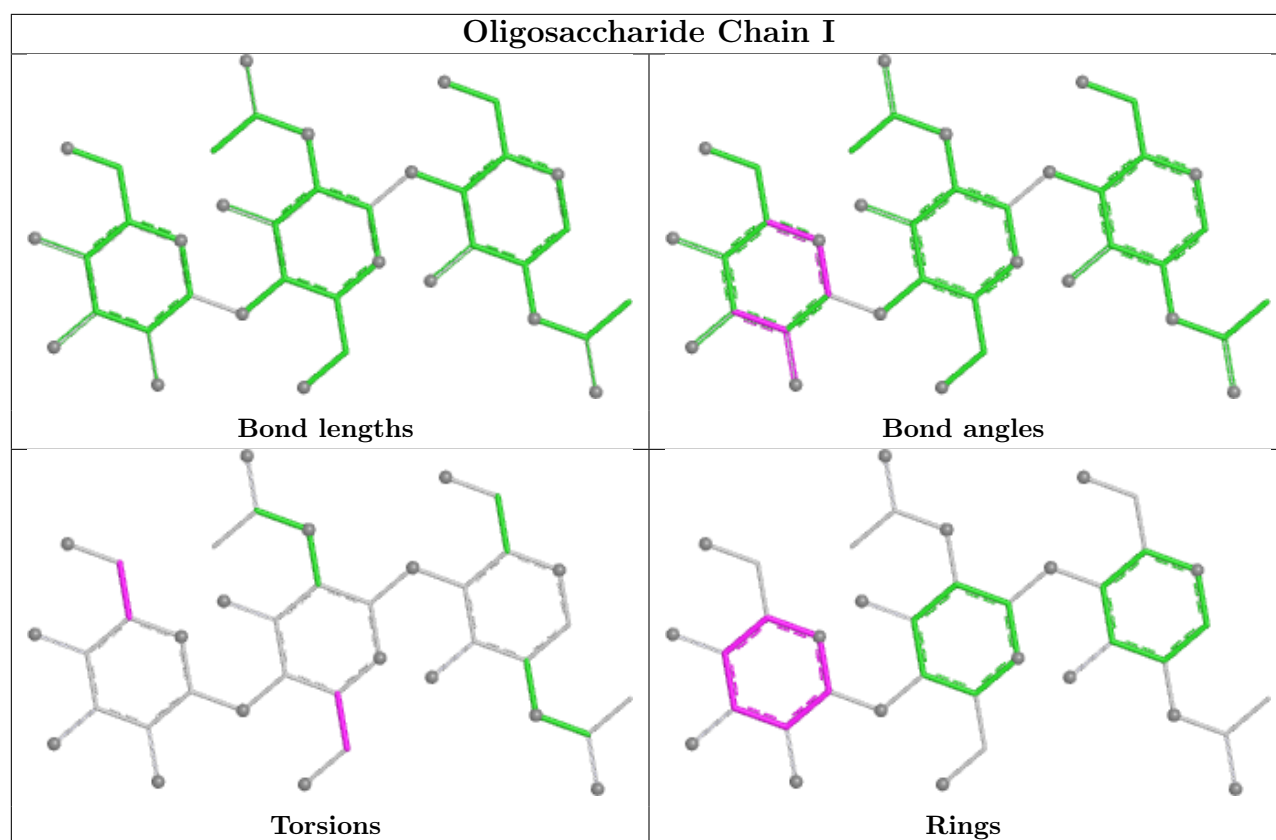
Mol	Chain	Res	Type	Atoms
2	L	3	MAN	C1-C2-C3-C4-C5-O5
2	W	3	MAN	C1-C2-C3-C4-C5-O5
2	d	3	MAN	C1-C2-C3-C4-C5-O5
2	J	3	MAN	C1-C2-C3-C4-C5-O5
2	V	3	MAN	C1-C2-C3-C4-C5-O5
2	a	3	MAN	C1-C2-C3-C4-C5-O5
2	M	3	MAN	C1-C2-C3-C4-C5-O5
2	f	3	MAN	C1-C2-C3-C4-C5-O5
2	Q	3	MAN	C1-C2-C3-C4-C5-O5
2	X	3	MAN	C1-C2-C3-C4-C5-O5
2	S	3	MAN	C1-C2-C3-C4-C5-O5
2	c	3	MAN	C1-C2-C3-C4-C5-O5
2	N	3	MAN	C1-C2-C3-C4-C5-O5
2	I	3	MAN	C1-C2-C3-C4-C5-O5

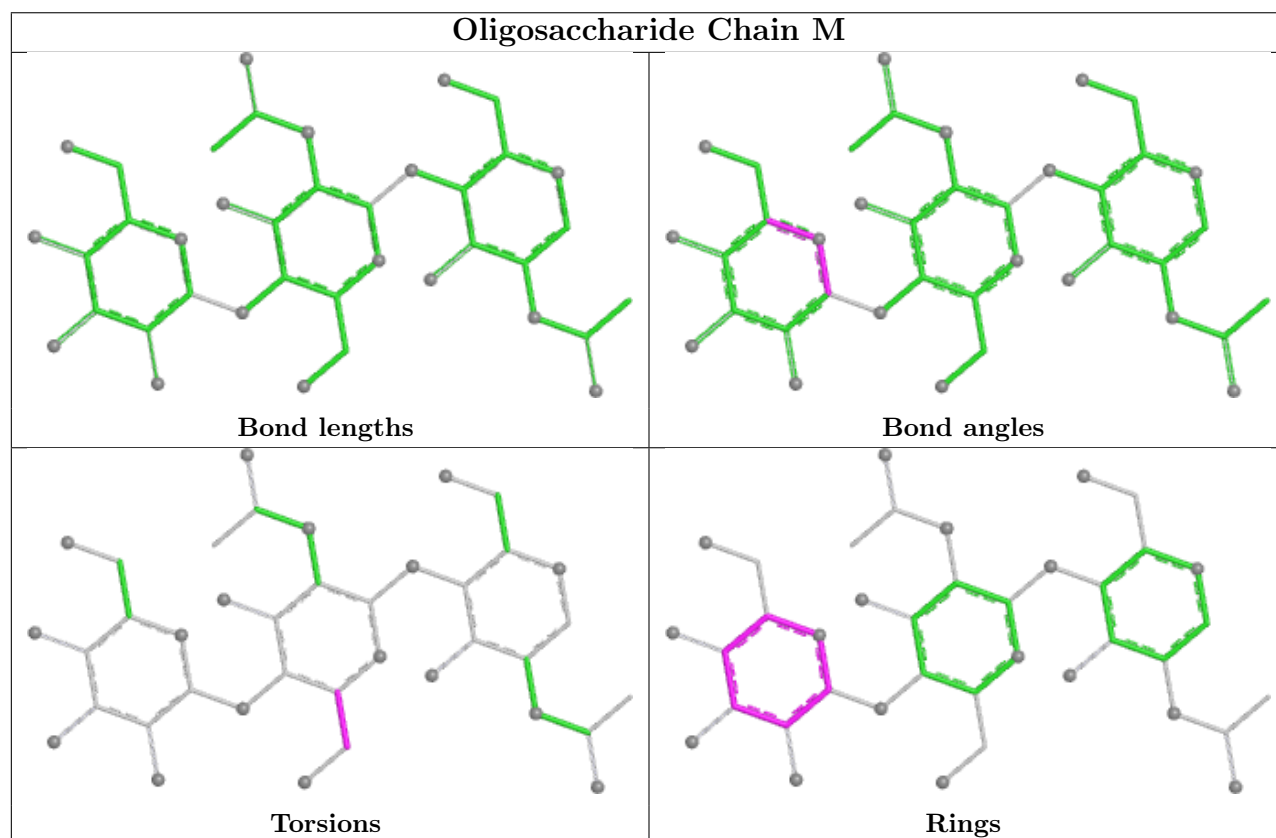
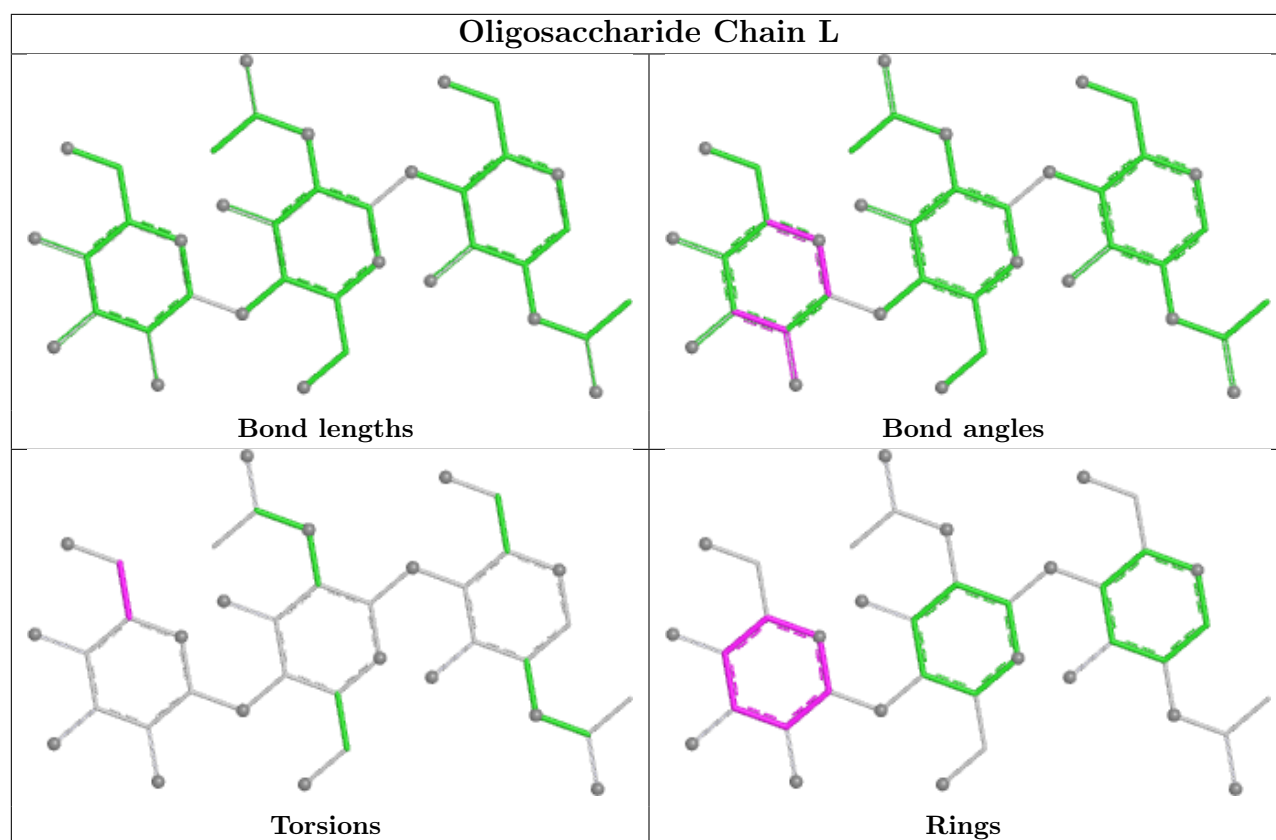
10 monomers are involved in 13 short contacts:

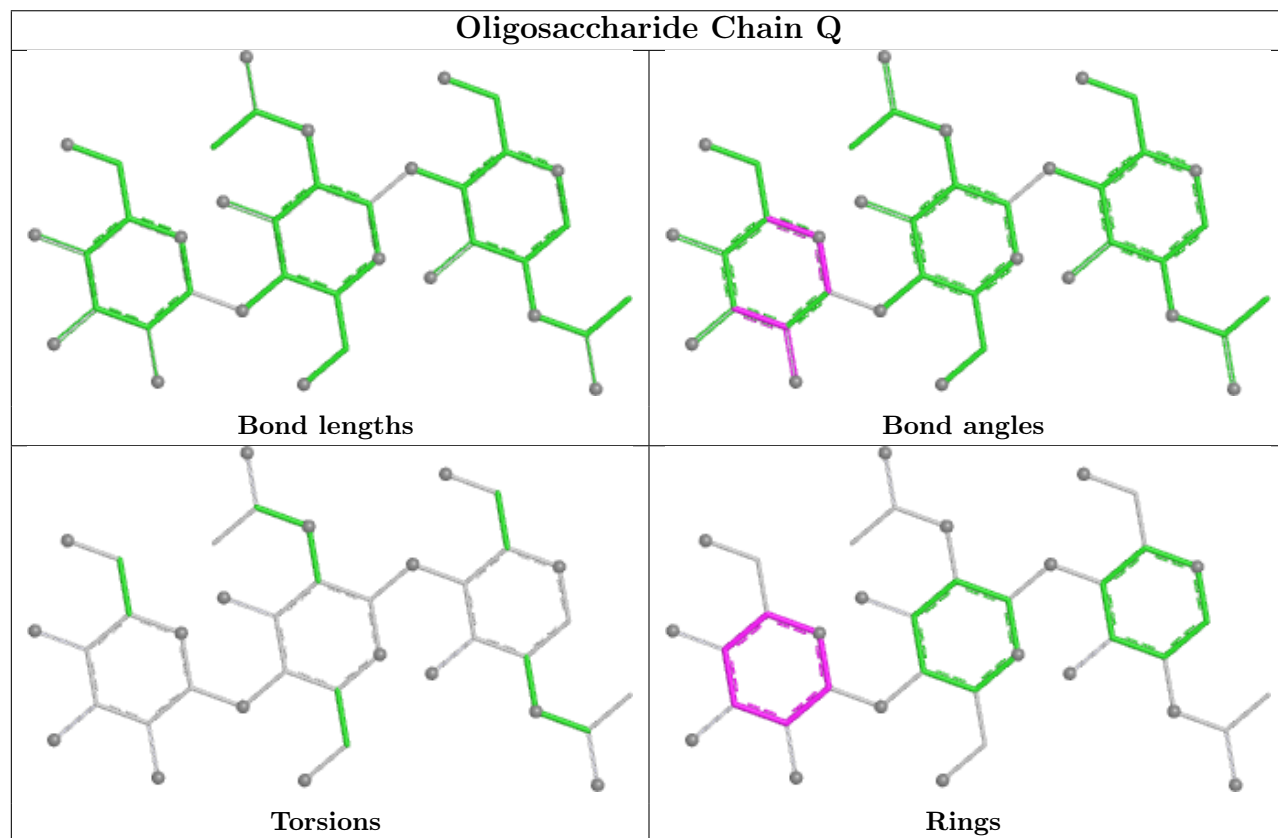
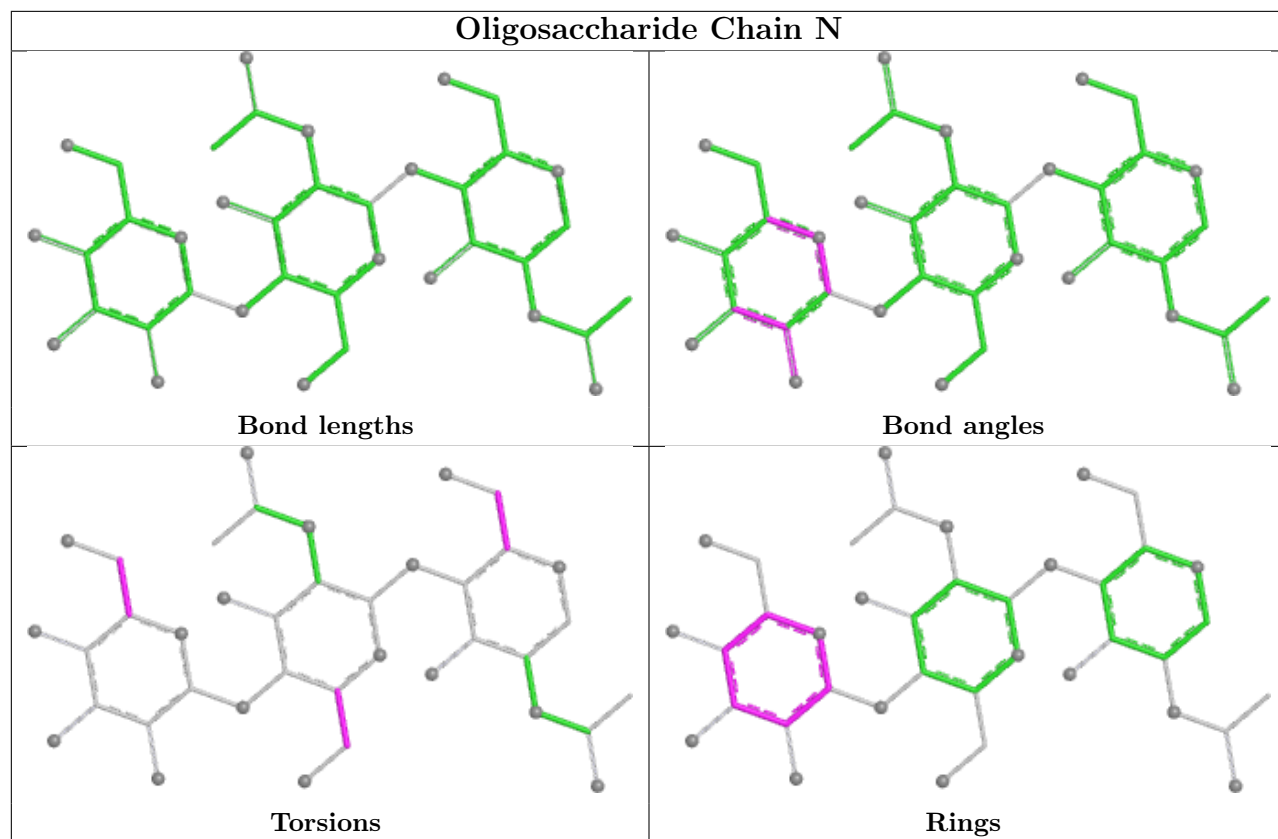
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	1	NAG	2	0
2	f	1	NAG	1	0
2	N	2	NAG	2	0
2	G	3	MAN	1	0
2	X	1	NAG	1	0
2	d	1	NAG	1	0
2	D	2	NAG	2	0
2	N	1	NAG	1	0
2	D	1	NAG	5	0
2	G	2	NAG	1	0

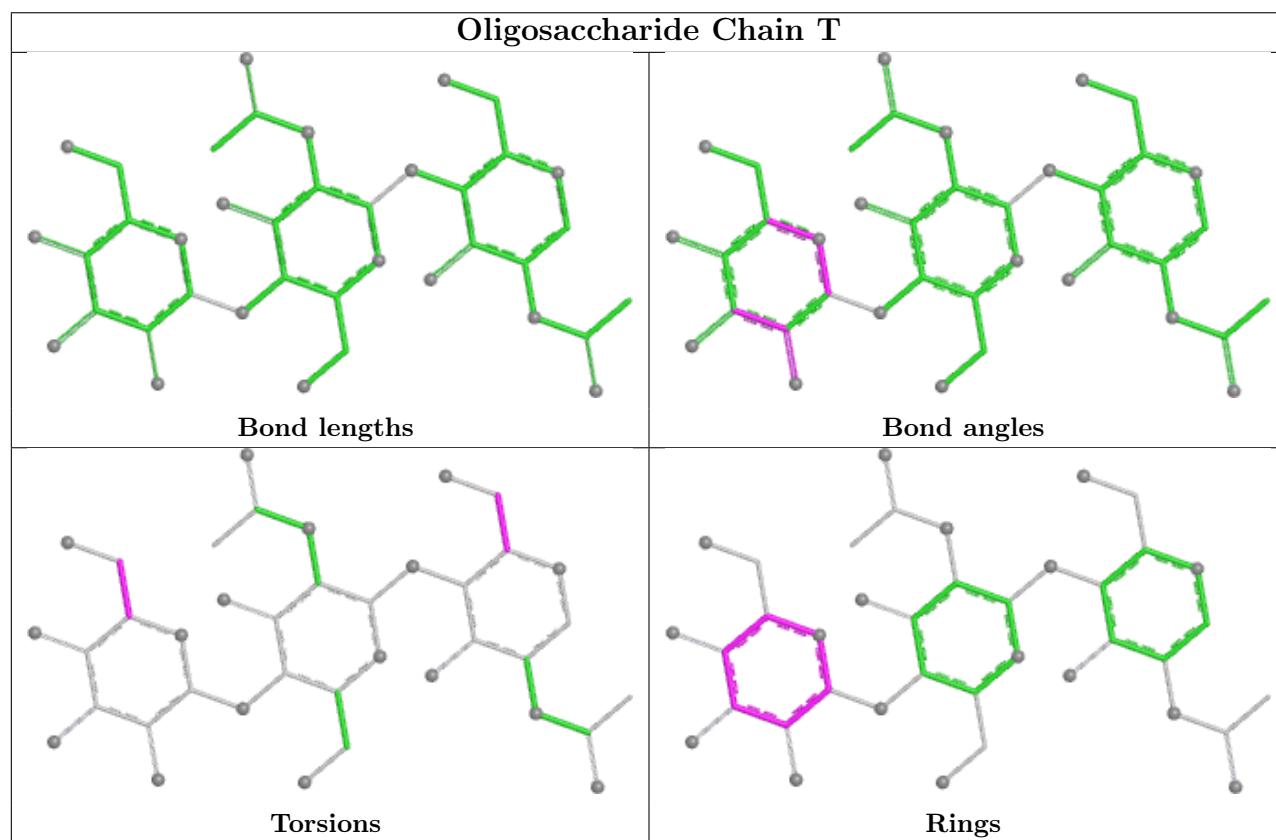
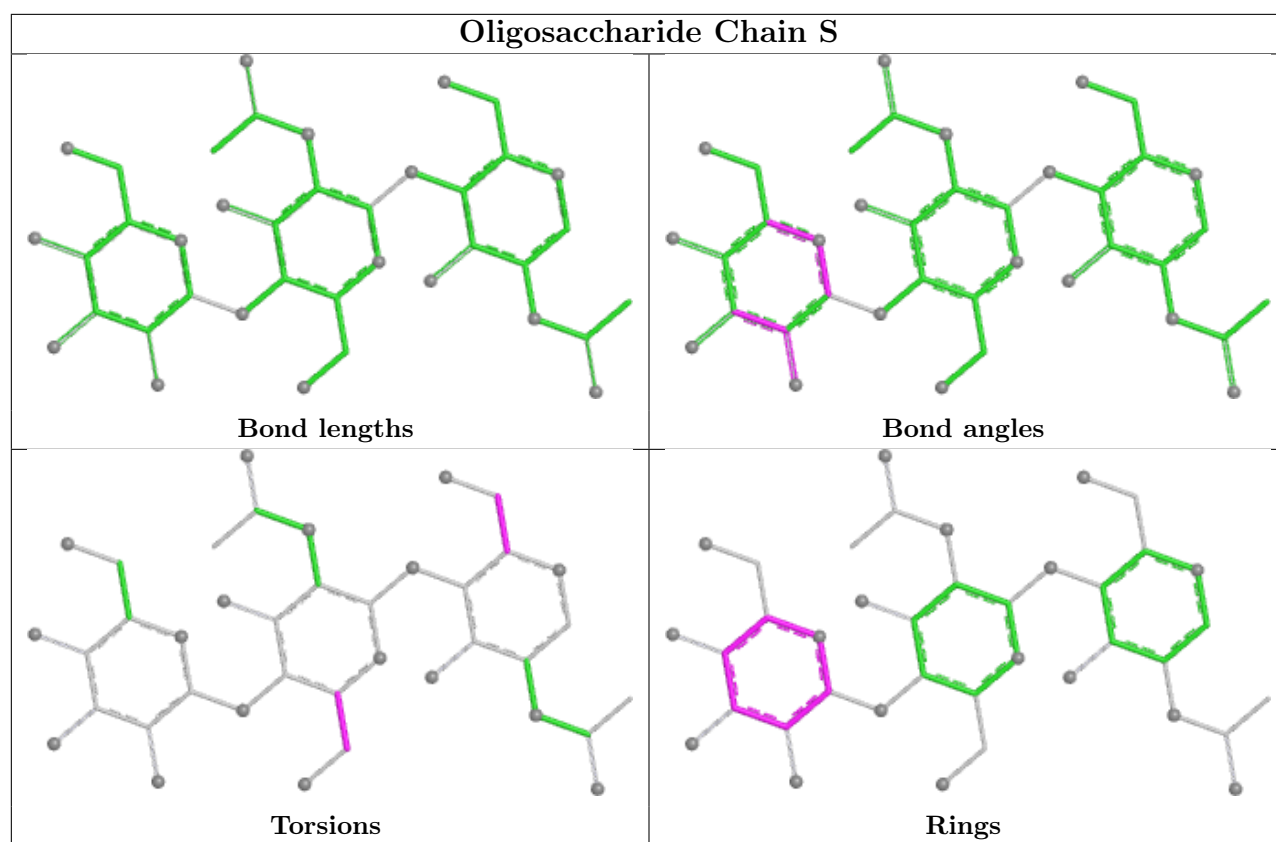
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

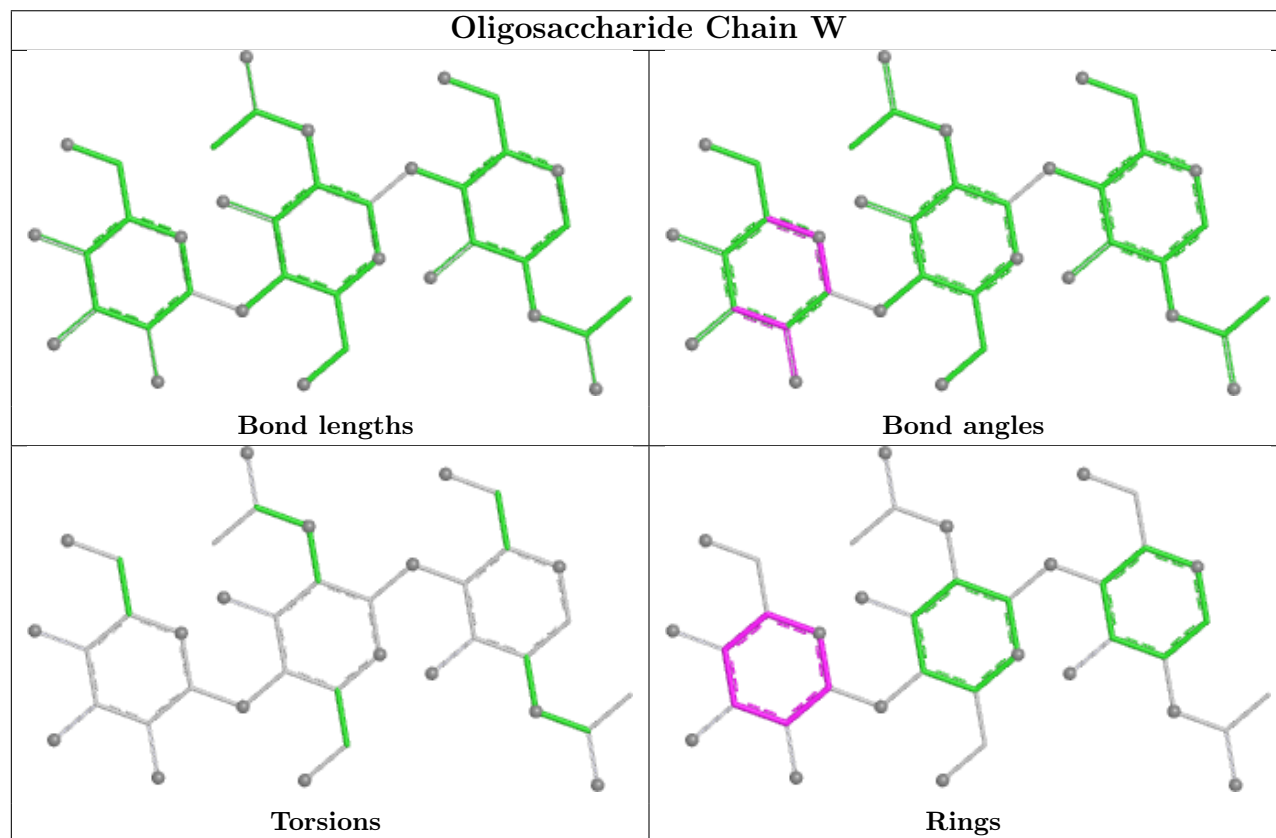
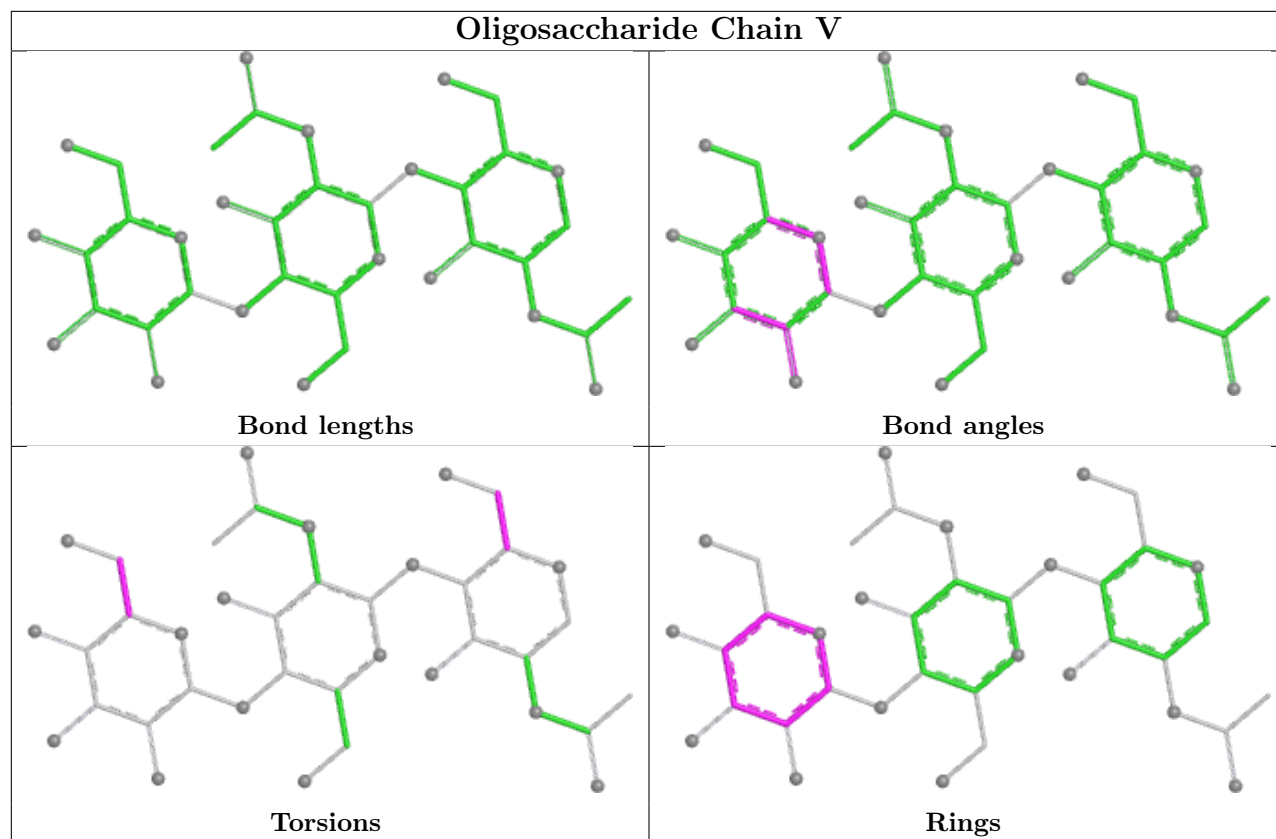


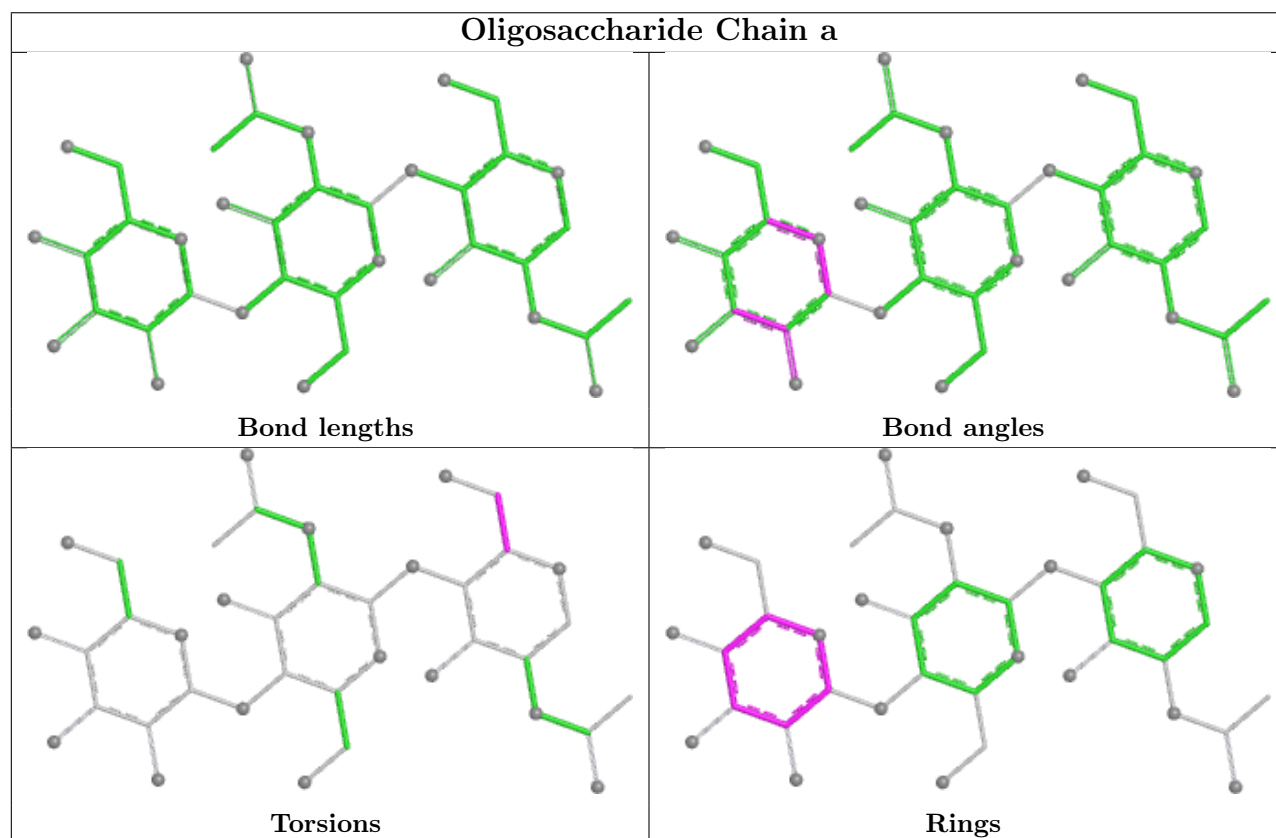
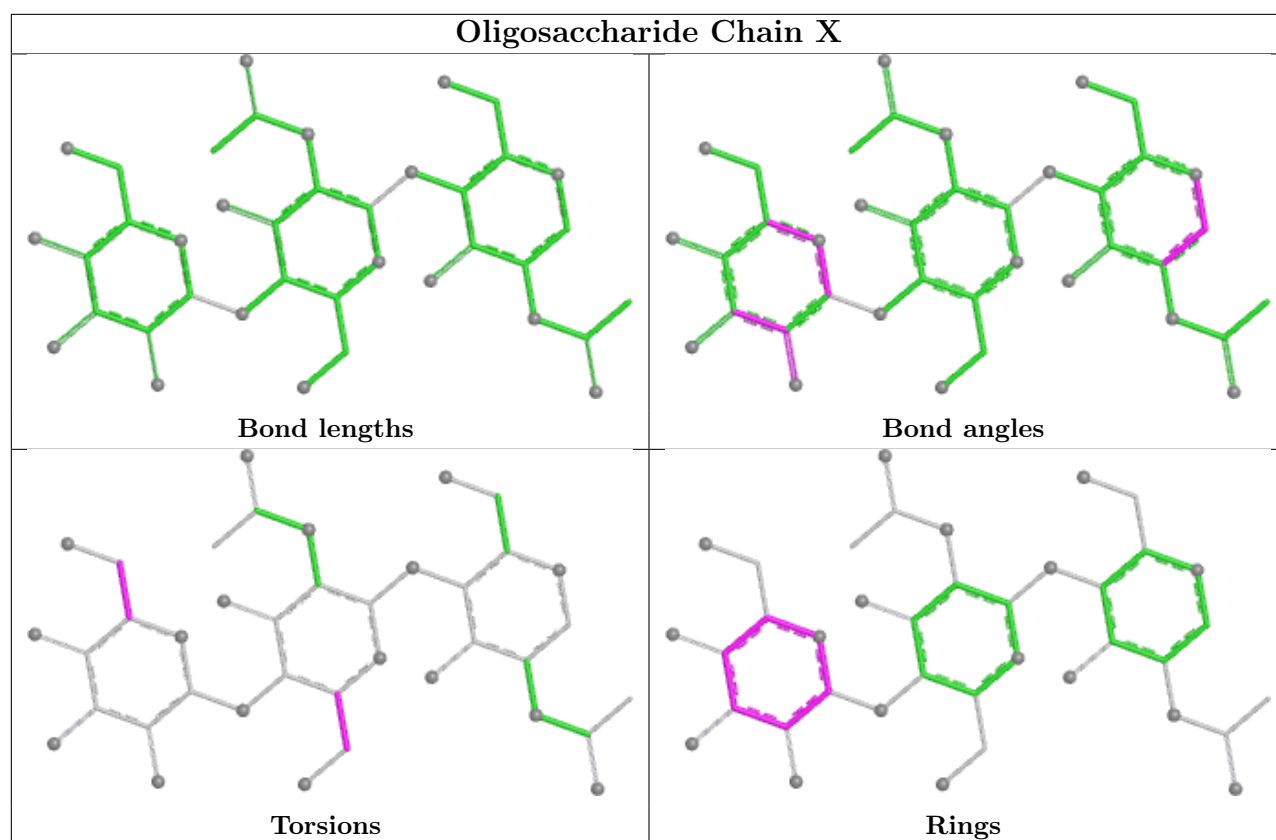


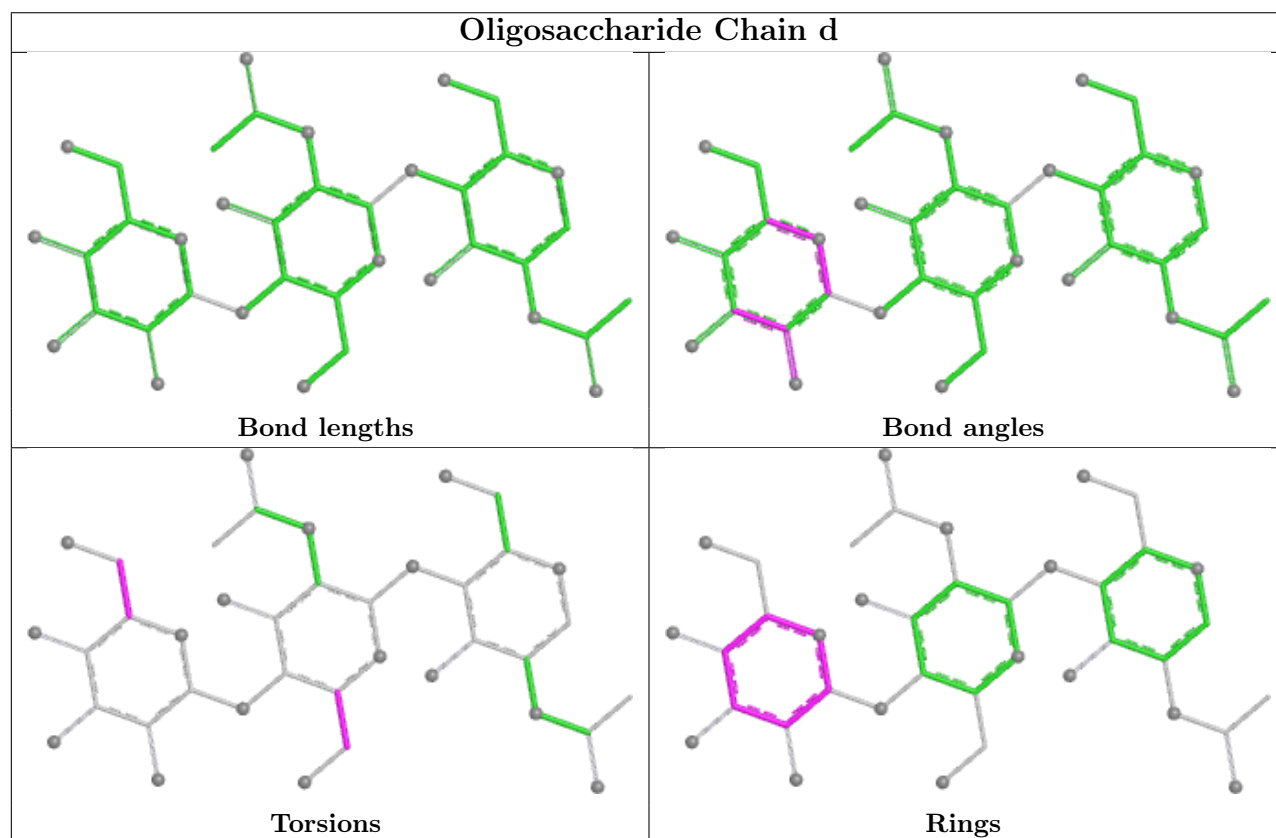
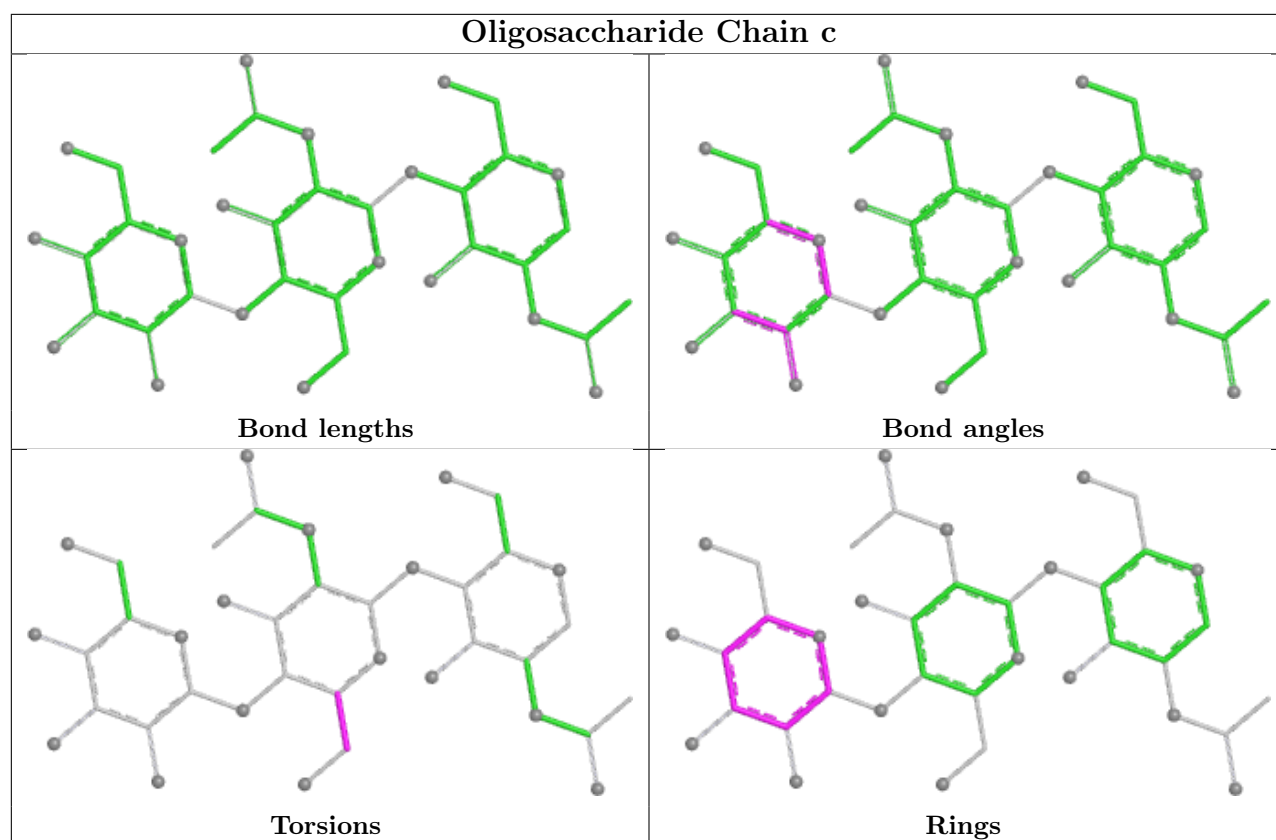


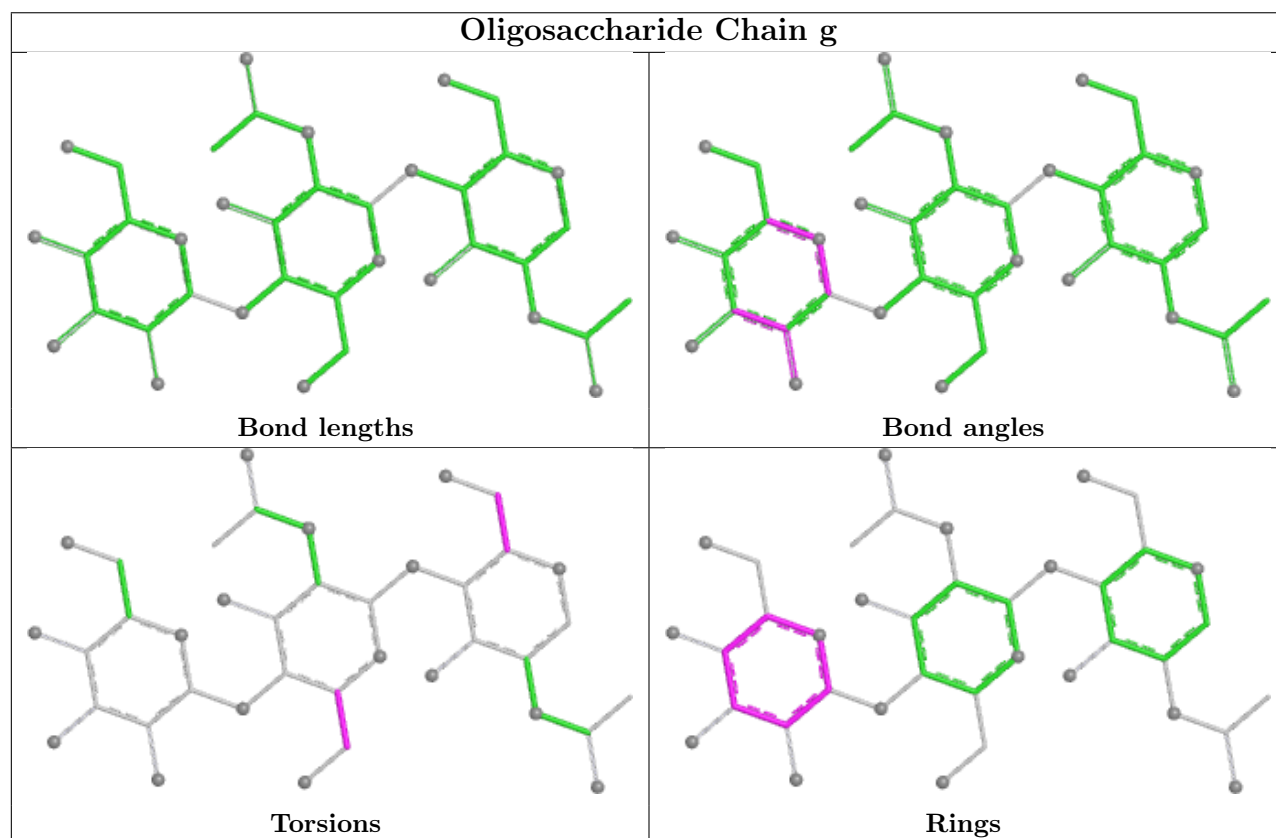
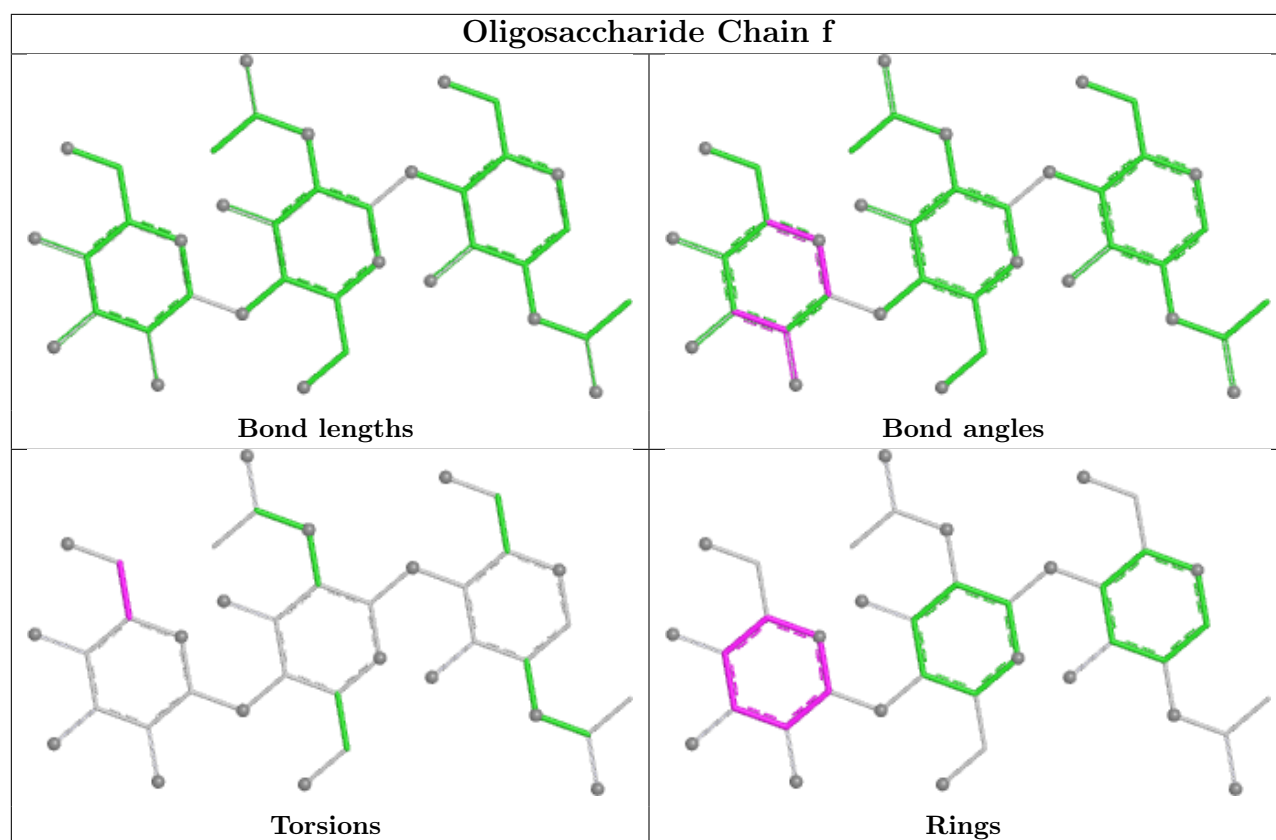


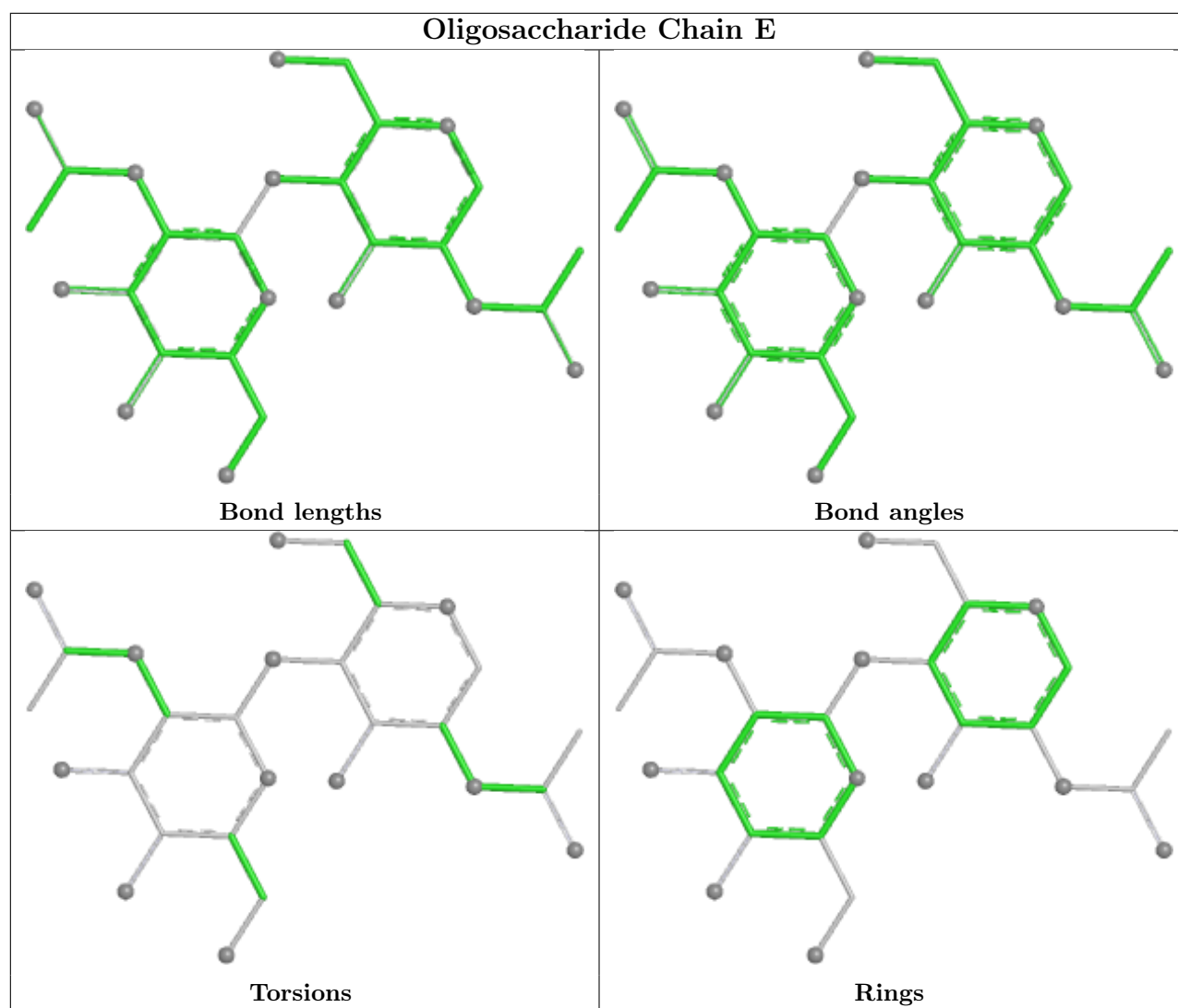


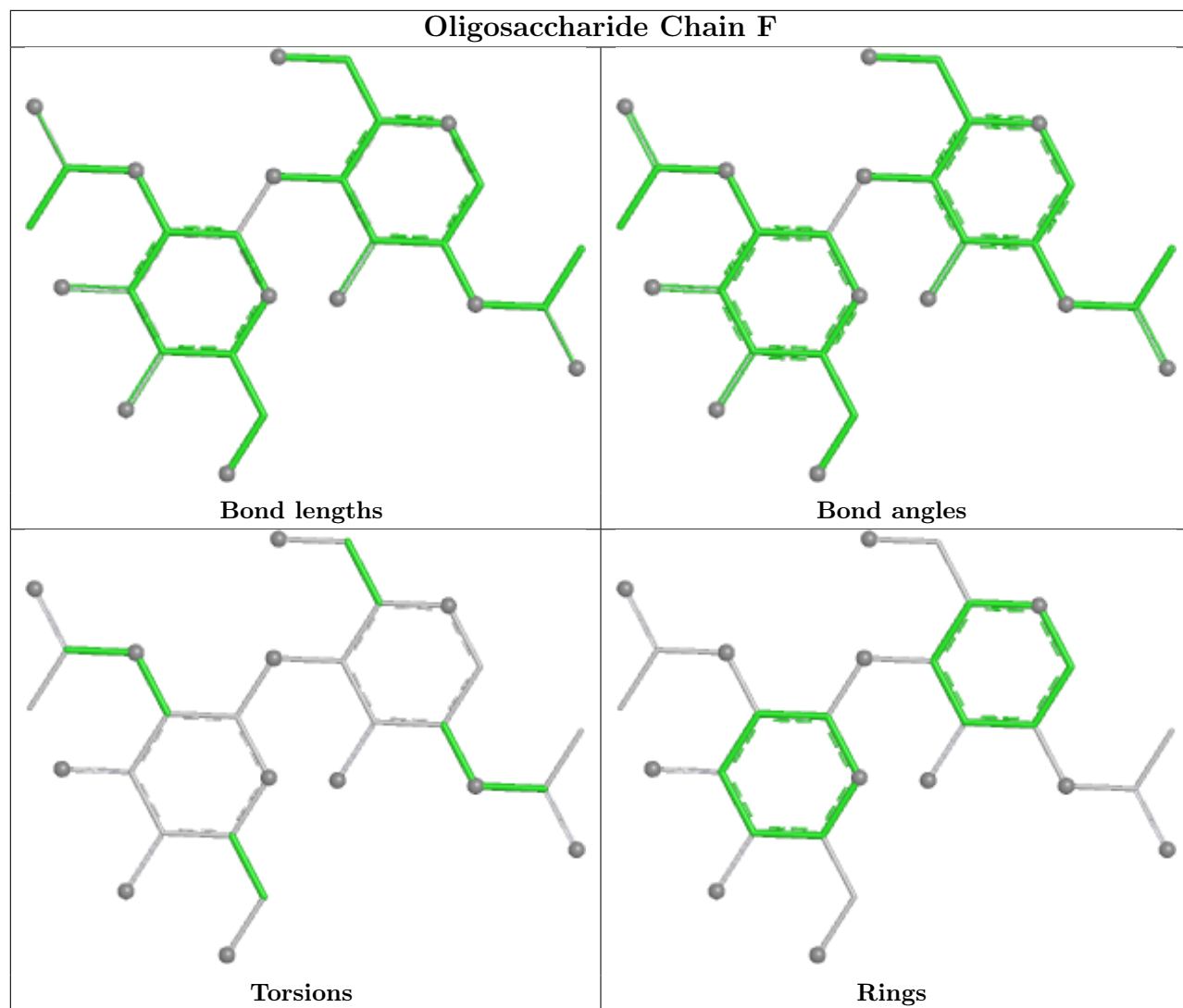


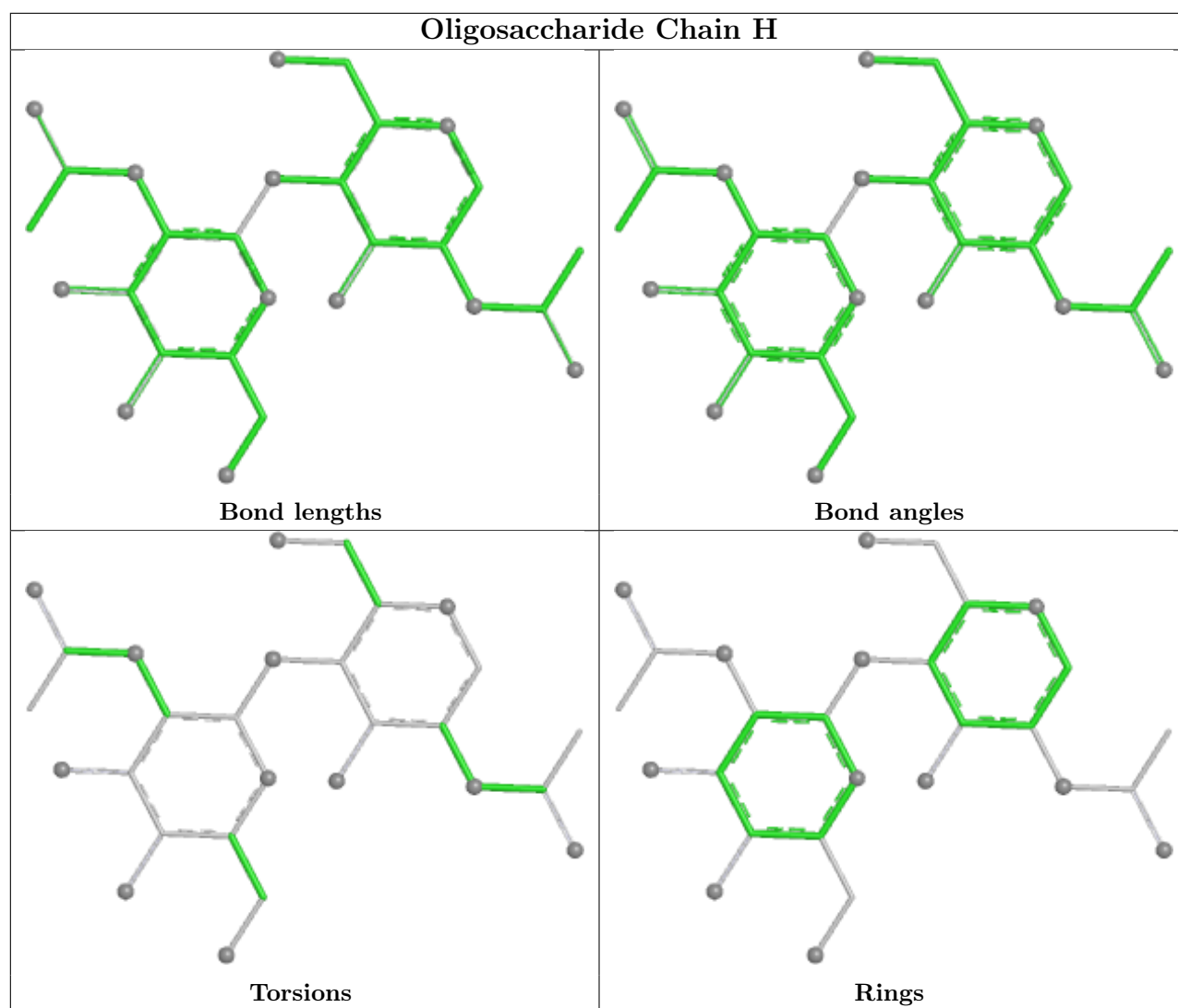


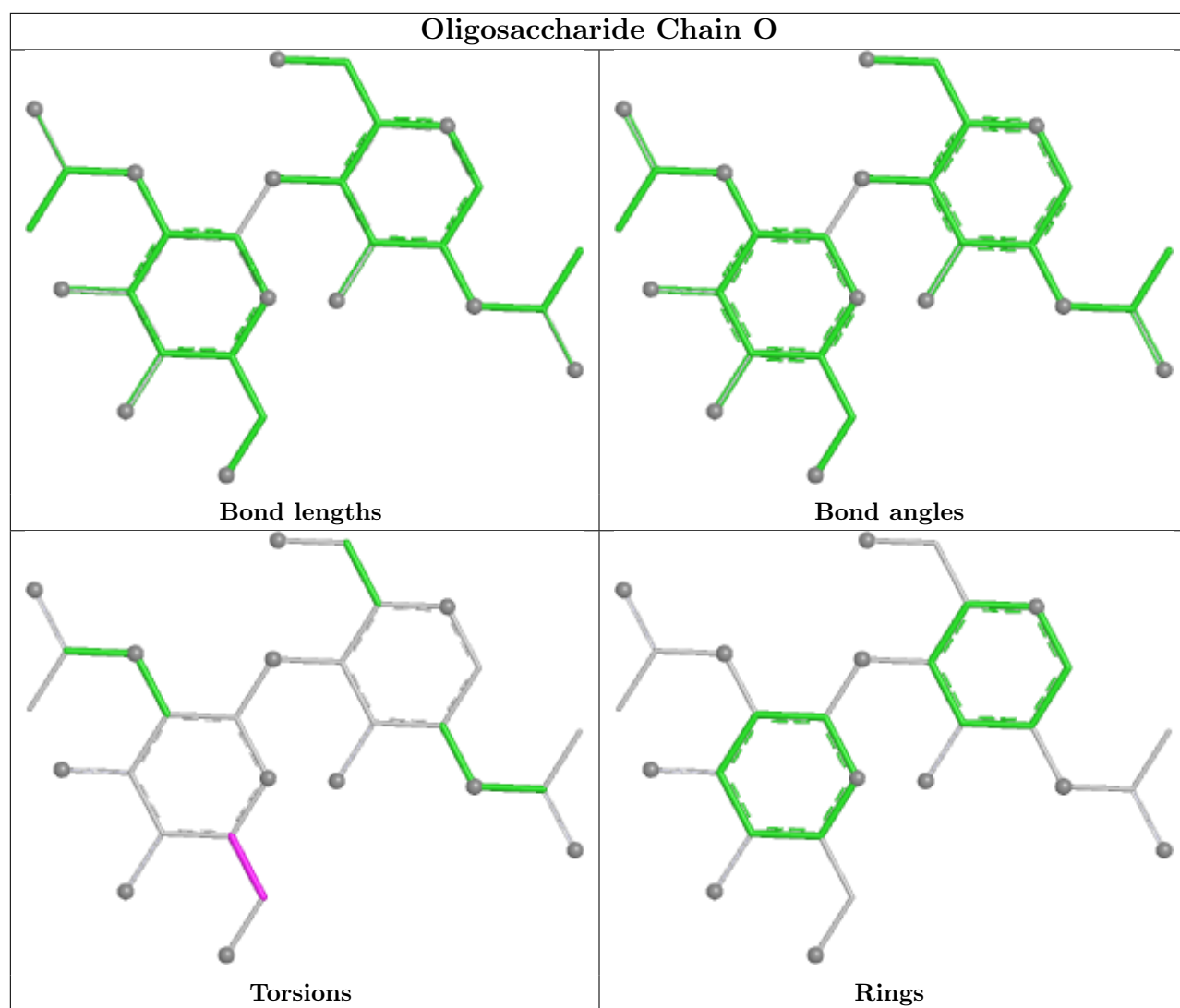


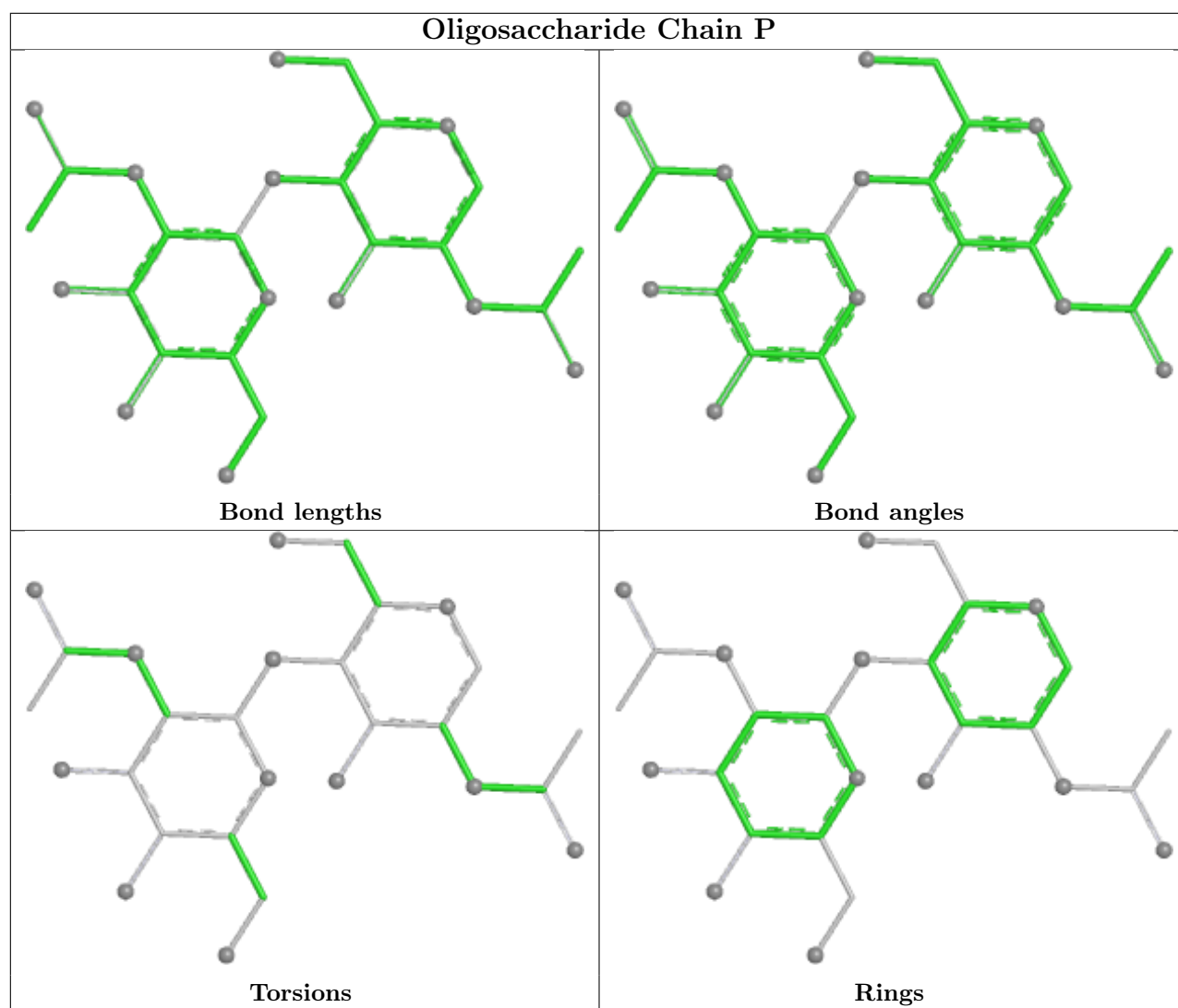


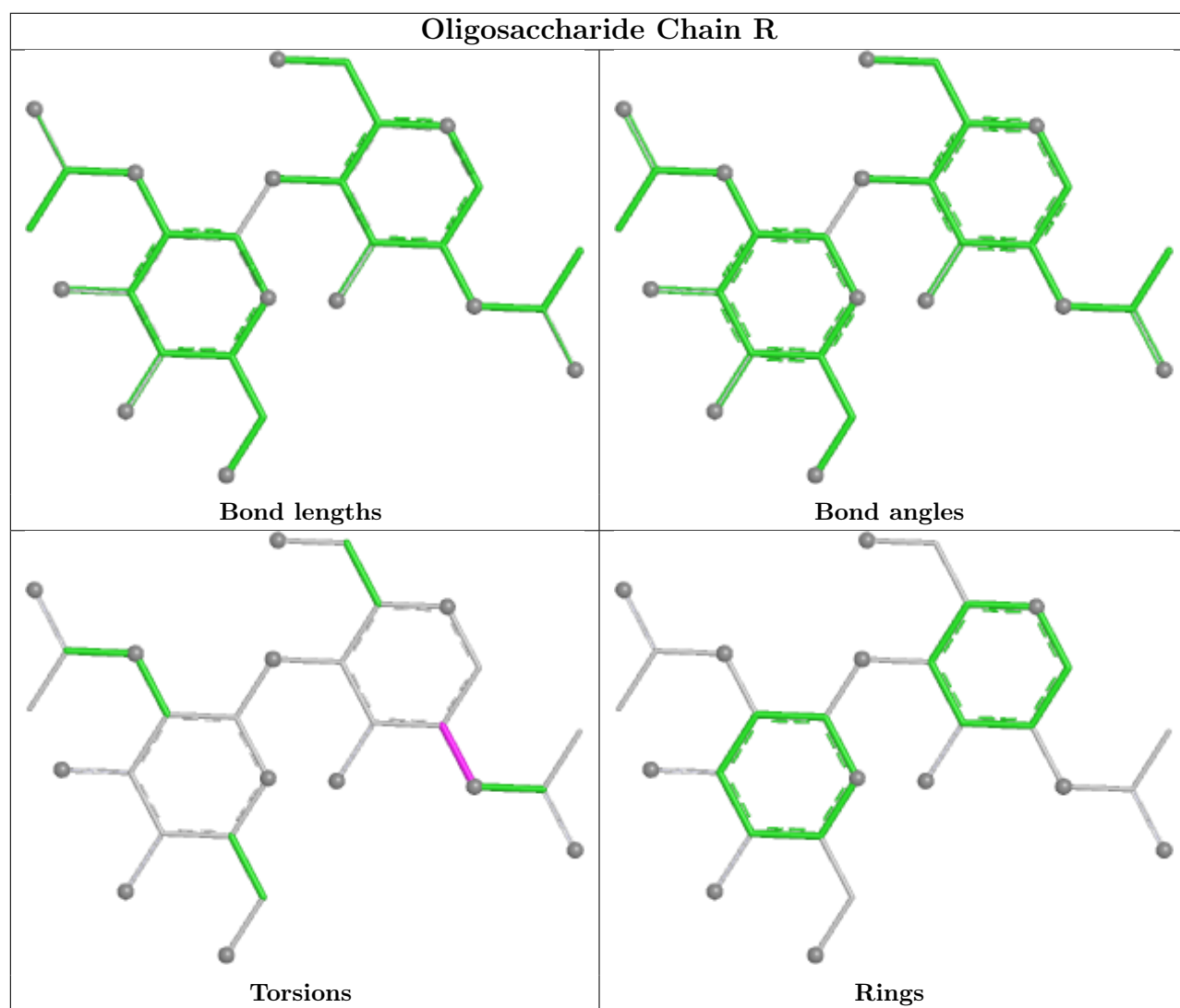


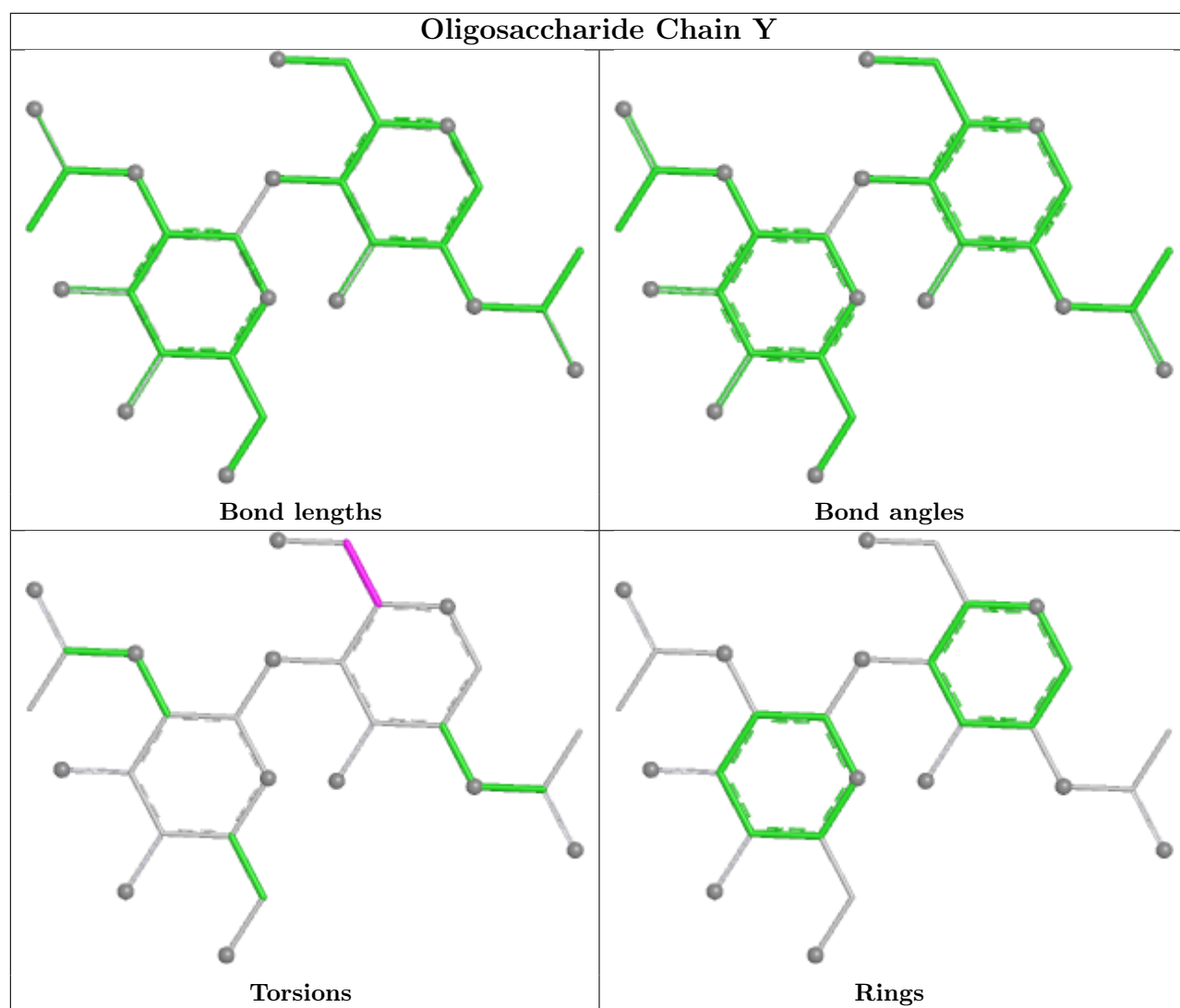


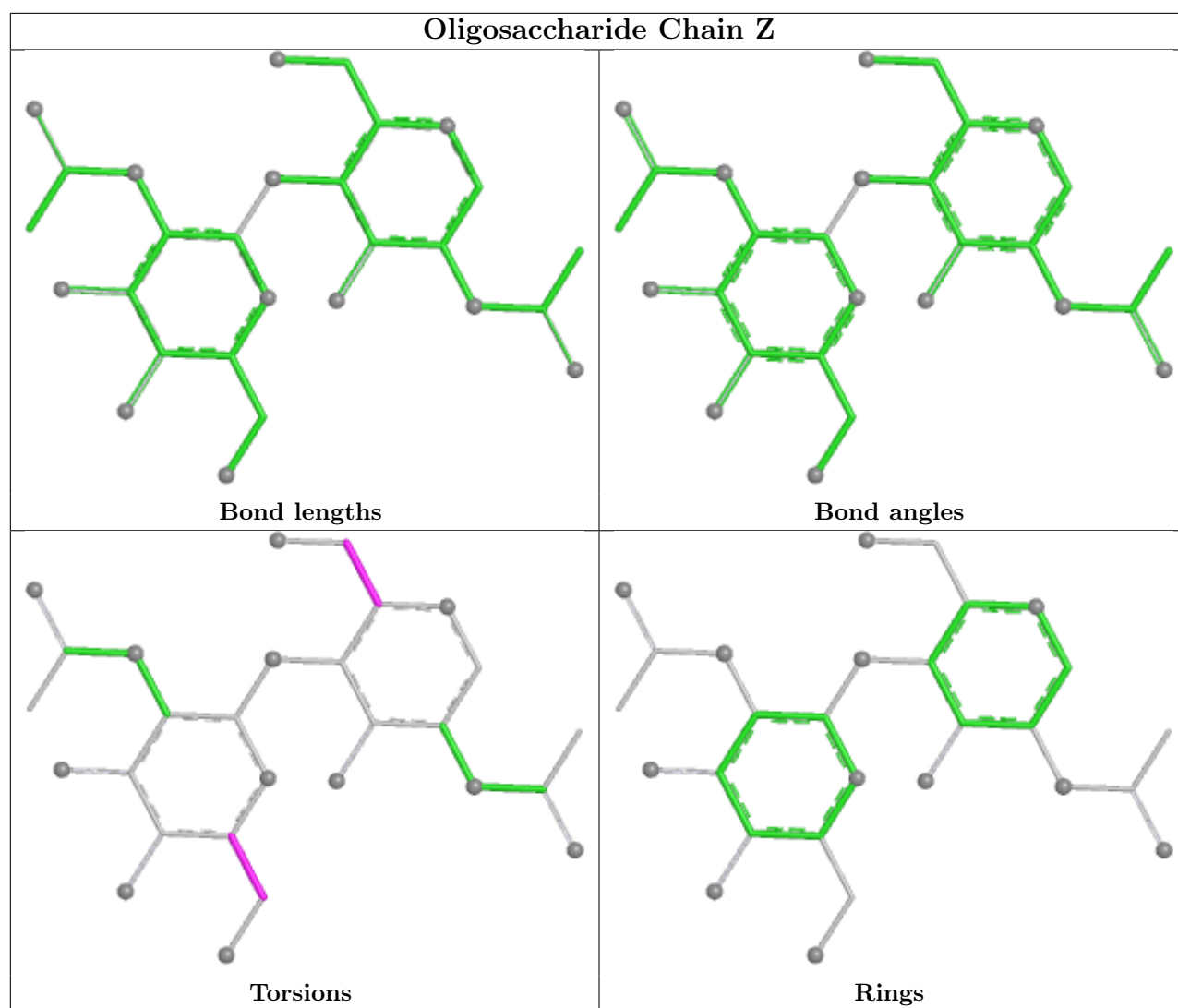


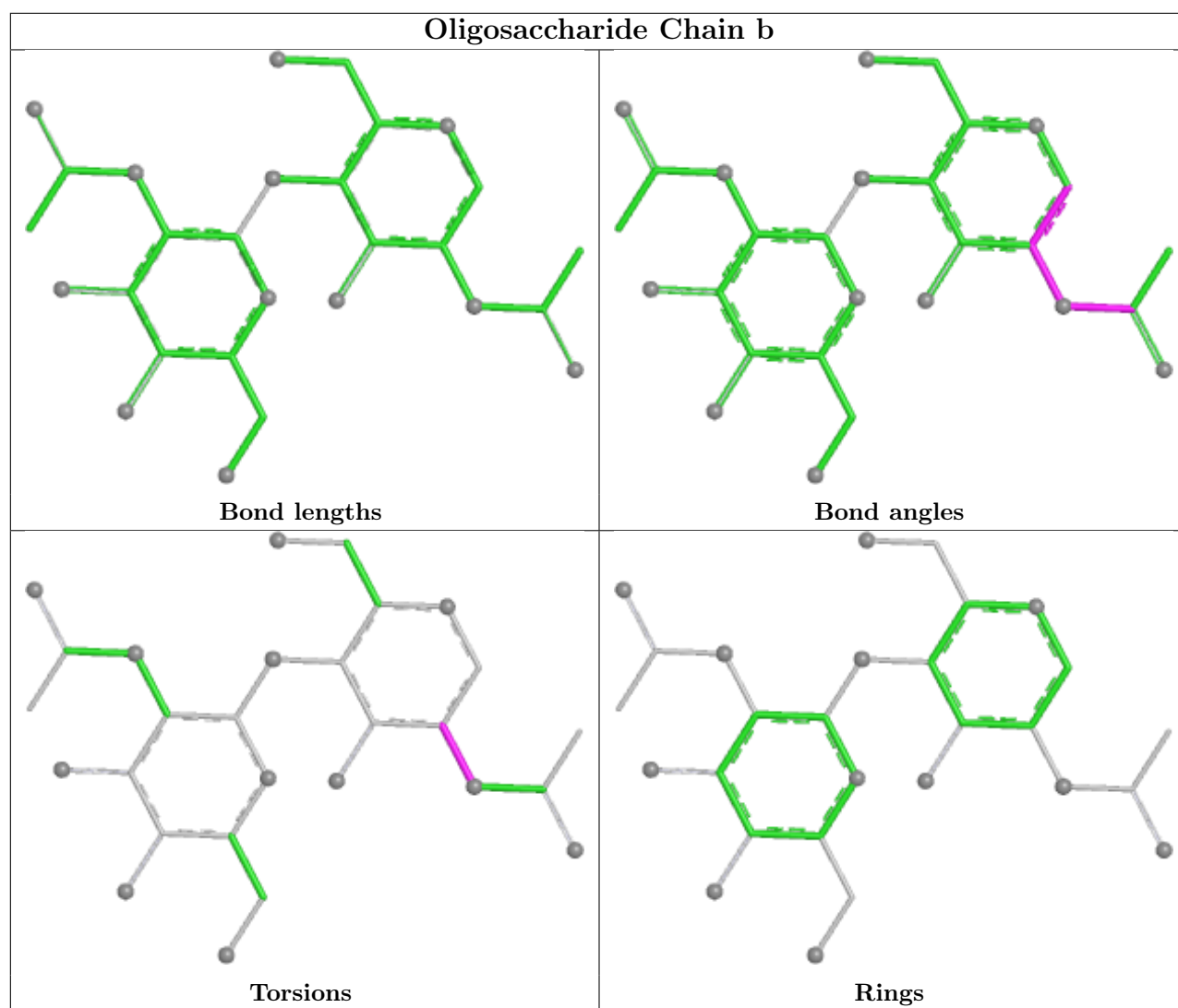


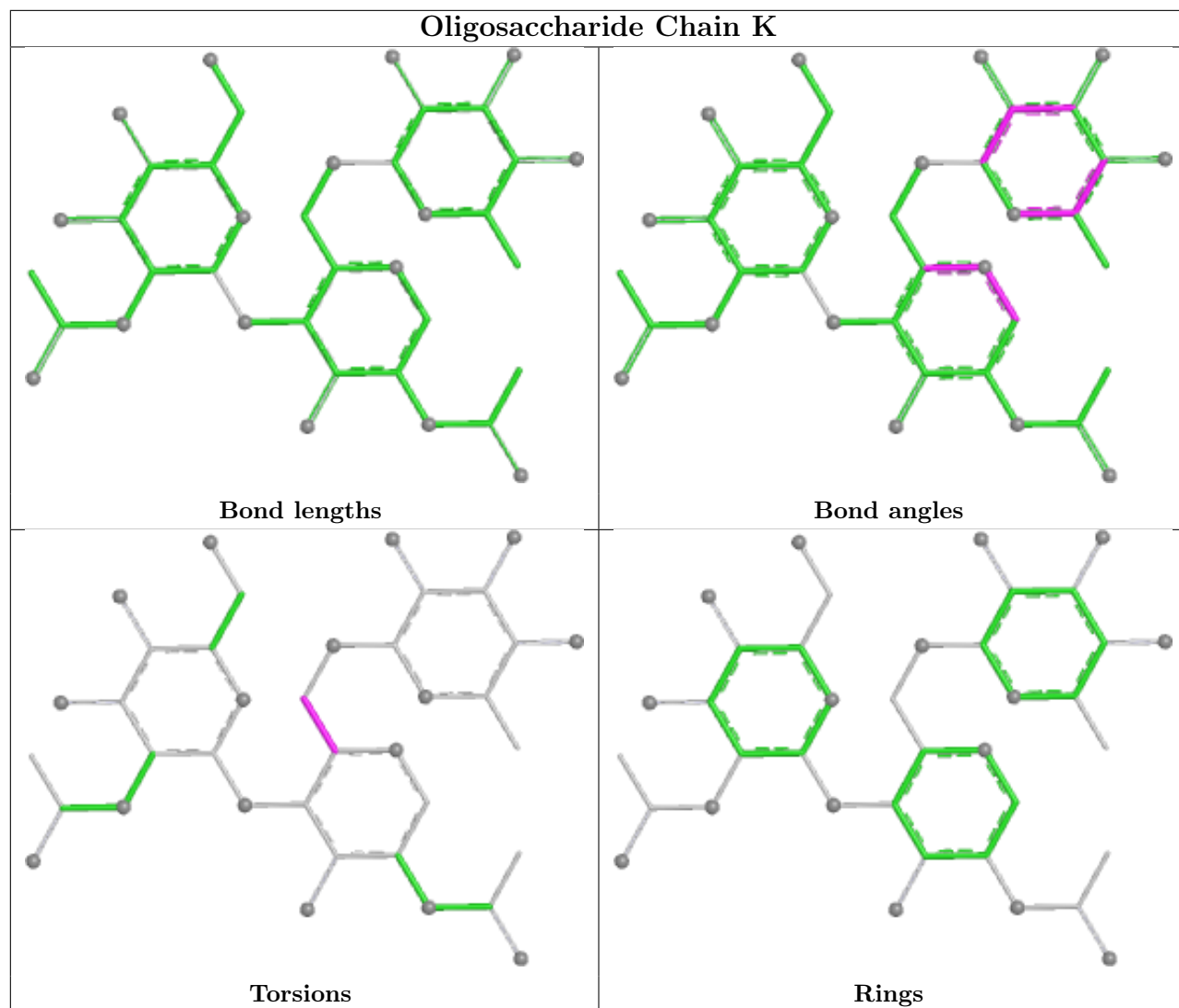


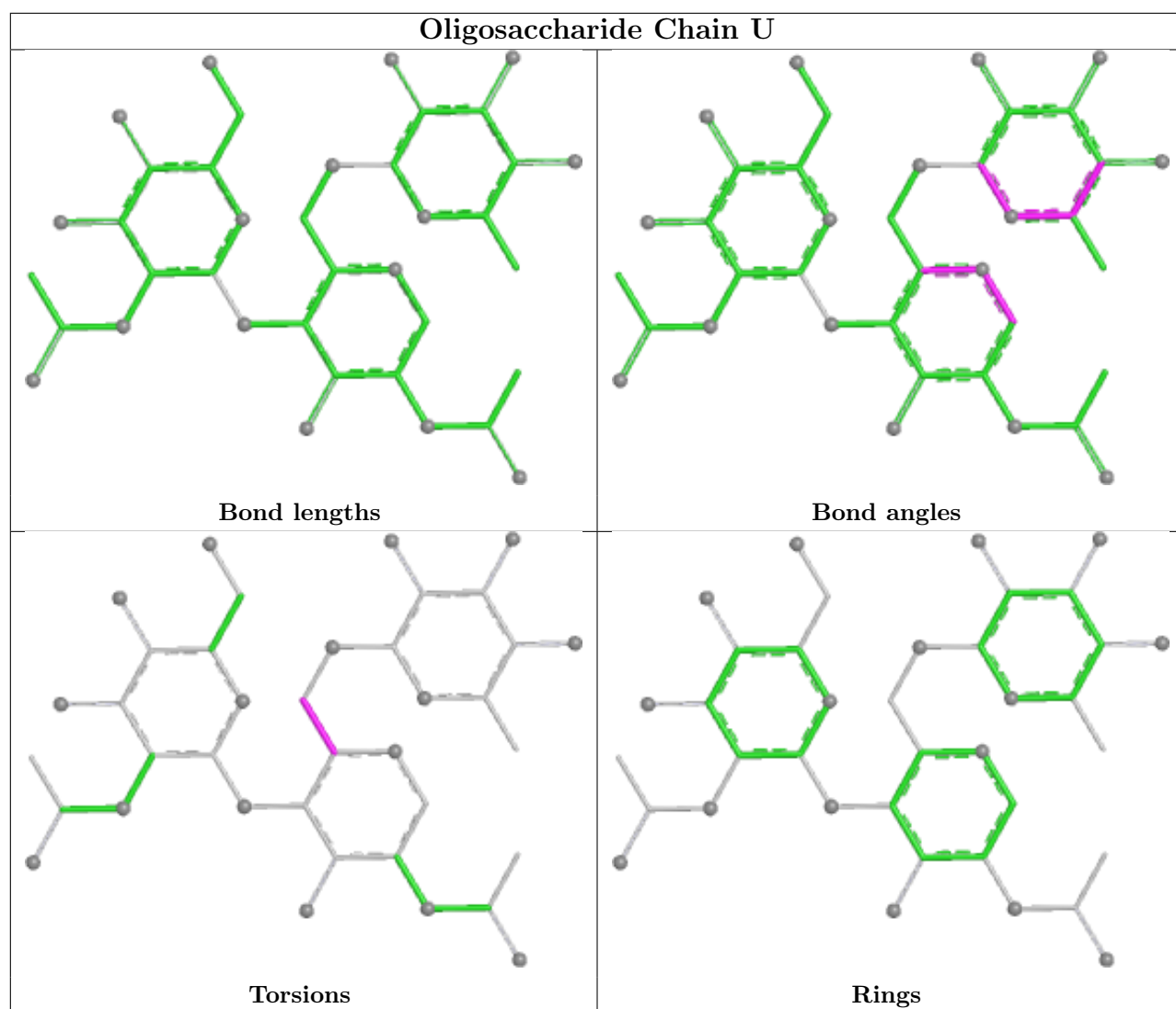


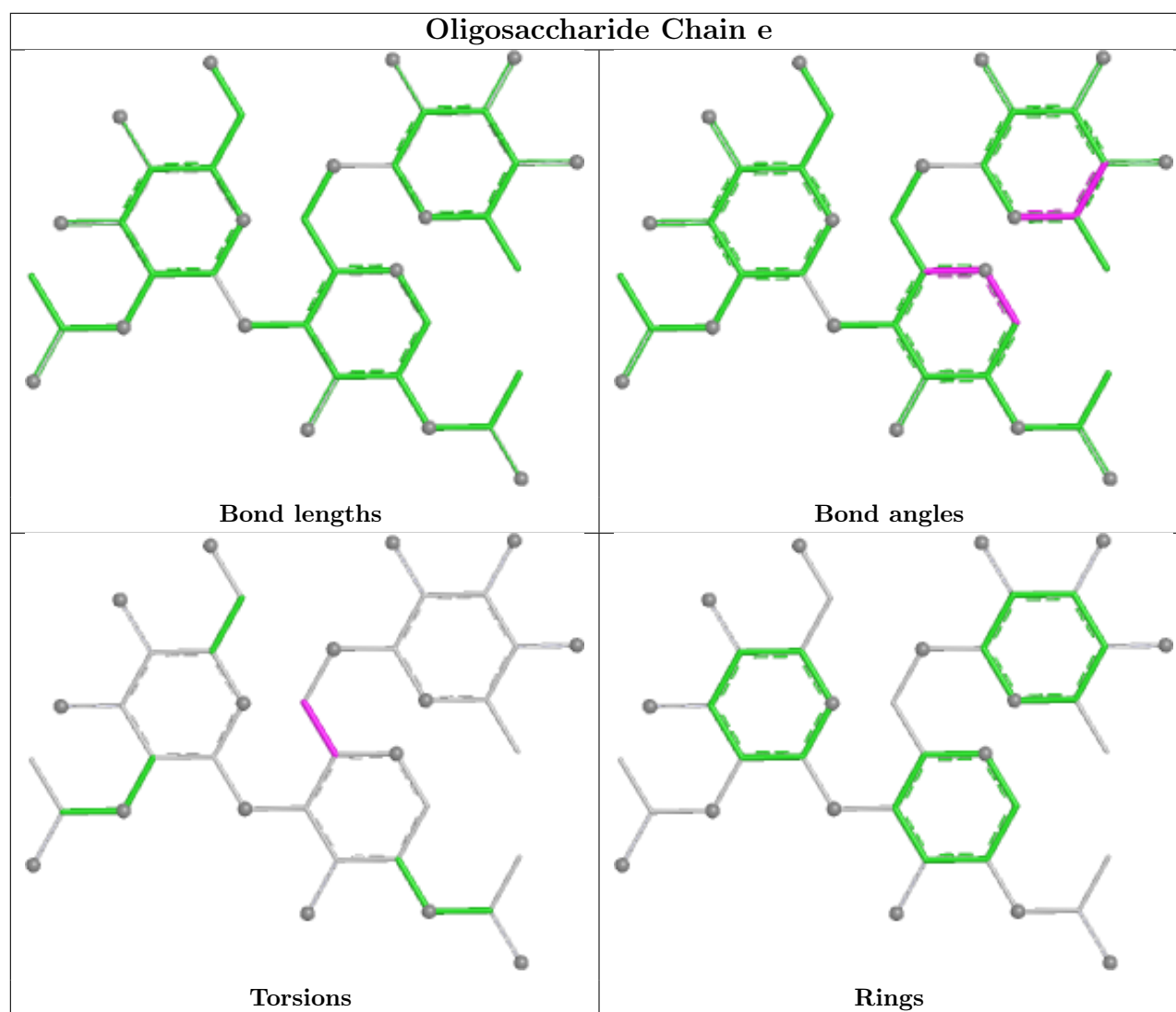












5.6 Ligand geometry [i](#)

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	$\# Z > 2$
5	NAG	B	1407	1	14,14,15	0.41	0	17,19,21	0.85	2 (11%)
5	NAG	B	1403	1	14,14,15	0.39	0	17,19,21	0.44	0
5	NAG	A	1407	1	14,14,15	0.40	0	17,19,21	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	C	1405	1	14,14,15	0.41	0	17,19,21	0.53	0
5	NAG	B	1401	1	14,14,15	0.41	0	17,19,21	0.59	0
5	NAG	A	1403	1	14,14,15	0.38	0	17,19,21	0.35	0
5	NAG	C	1407	1	14,14,15	0.40	0	17,19,21	0.35	0
5	NAG	A	1404	1	14,14,15	0.41	0	17,19,21	0.36	0
5	NAG	B	1404	1	14,14,15	0.40	0	17,19,21	0.34	0
5	NAG	C	1404	1	14,14,15	0.40	0	17,19,21	0.35	0
5	NAG	B	1405	1	14,14,15	0.42	0	17,19,21	0.38	0
5	NAG	A	1402	1	14,14,15	0.40	0	17,19,21	0.43	0
5	NAG	A	1405	1	14,14,15	0.40	0	17,19,21	0.50	0
5	NAG	C	1402	1	14,14,15	0.40	0	17,19,21	0.48	0
5	NAG	B	1402	1	14,14,15	0.41	0	17,19,21	0.44	0
5	NAG	C	1406	1	14,14,15	0.42	0	17,19,21	0.39	0
5	NAG	A	1401	1	14,14,15	0.29	0	17,19,21	0.47	0
5	NAG	C	1401	1	14,14,15	0.28	0	17,19,21	0.47	0
5	NAG	A	1406	1	14,14,15	0.39	0	17,19,21	0.36	0
5	NAG	C	1403	1	14,14,15	0.40	0	17,19,21	0.50	0
5	NAG	B	1406	1	14,14,15	0.41	0	17,19,21	0.40	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	B	1407	1	-	4/6/23/26	0/1/1/1
5	NAG	B	1403	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1407	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1405	1	-	1/6/23/26	0/1/1/1
5	NAG	B	1401	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1403	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1407	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1404	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1404	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1404	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1405	1	-	3/6/23/26	0/1/1/1
5	NAG	A	1402	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1405	1	-	1/6/23/26	0/1/1/1
5	NAG	C	1402	1	-	1/6/23/26	0/1/1/1
5	NAG	B	1402	1	-	3/6/23/26	0/1/1/1
5	NAG	C	1406	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1401	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	C	1401	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1406	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1403	1	-	1/6/23/26	0/1/1/1
5	NAG	B	1406	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1407	NAG	C1-C2-N2	2.21	113.92	110.43
5	B	1407	NAG	C2-N2-C7	2.19	125.84	122.90

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1407	NAG	C1-C2-N2-C7
5	A	1401	NAG	O5-C5-C6-O6
5	B	1402	NAG	C8-C7-N2-C2
5	B	1402	NAG	O7-C7-N2-C2
5	A	1401	NAG	C4-C5-C6-O6
5	B	1407	NAG	C8-C7-N2-C2
5	B	1407	NAG	O7-C7-N2-C2
5	B	1407	NAG	C3-C2-N2-C7
5	B	1405	NAG	C8-C7-N2-C2
5	B	1405	NAG	O7-C7-N2-C2
5	A	1402	NAG	C3-C2-N2-C7
5	A	1405	NAG	C3-C2-N2-C7
5	B	1401	NAG	C3-C2-N2-C7
5	B	1402	NAG	C3-C2-N2-C7
5	B	1405	NAG	C3-C2-N2-C7
5	C	1402	NAG	C3-C2-N2-C7
5	C	1403	NAG	C3-C2-N2-C7
5	C	1405	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 4 short contacts:

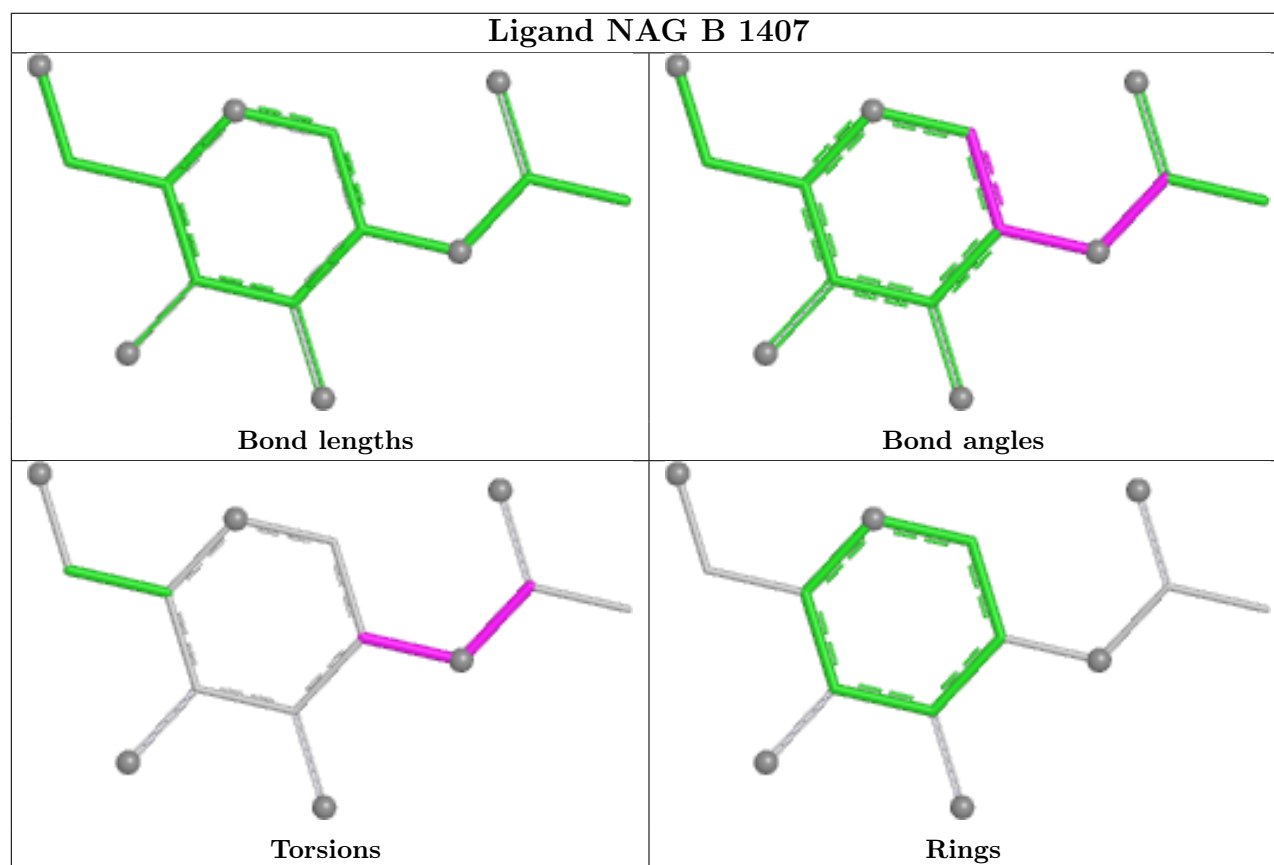
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1403	NAG	2	0

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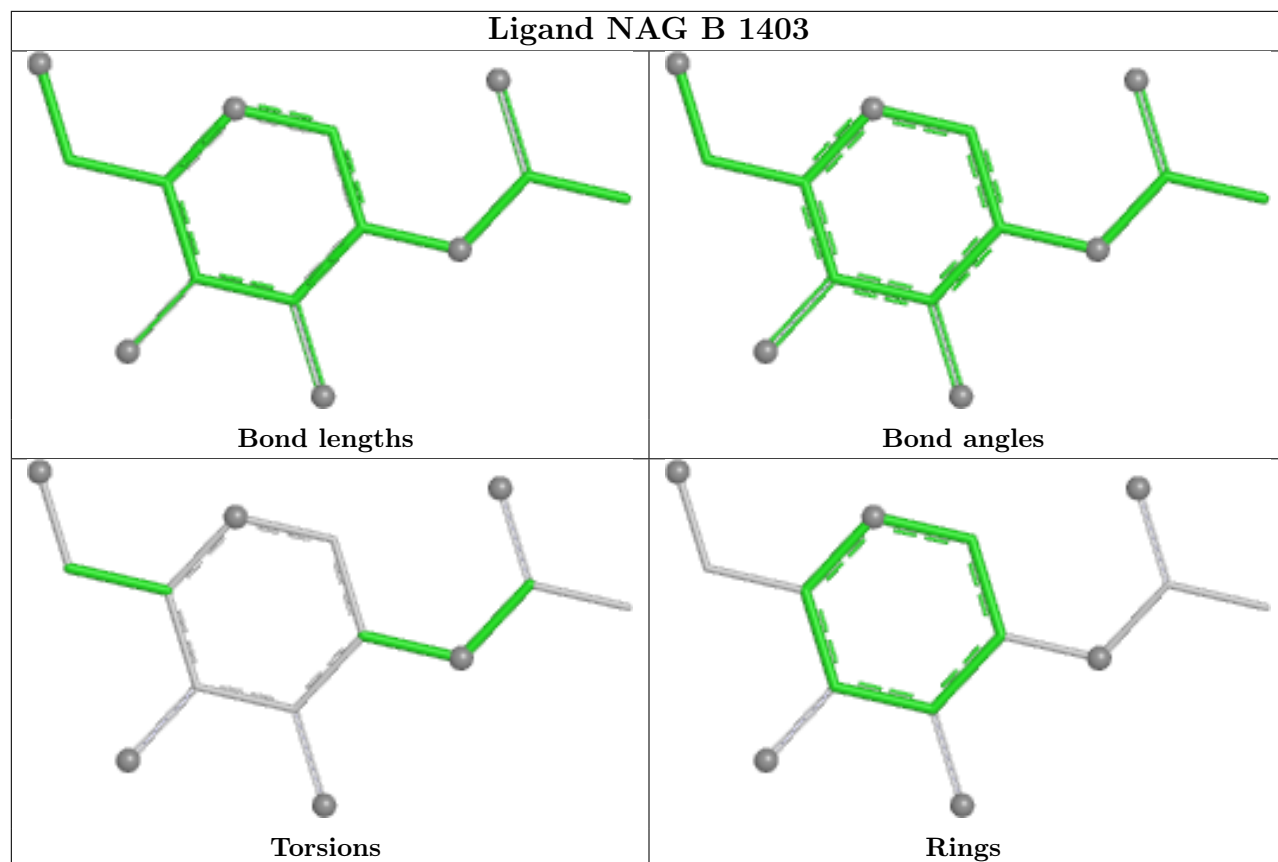
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1403	NAG	1	0
5	A	1401	NAG	1	0

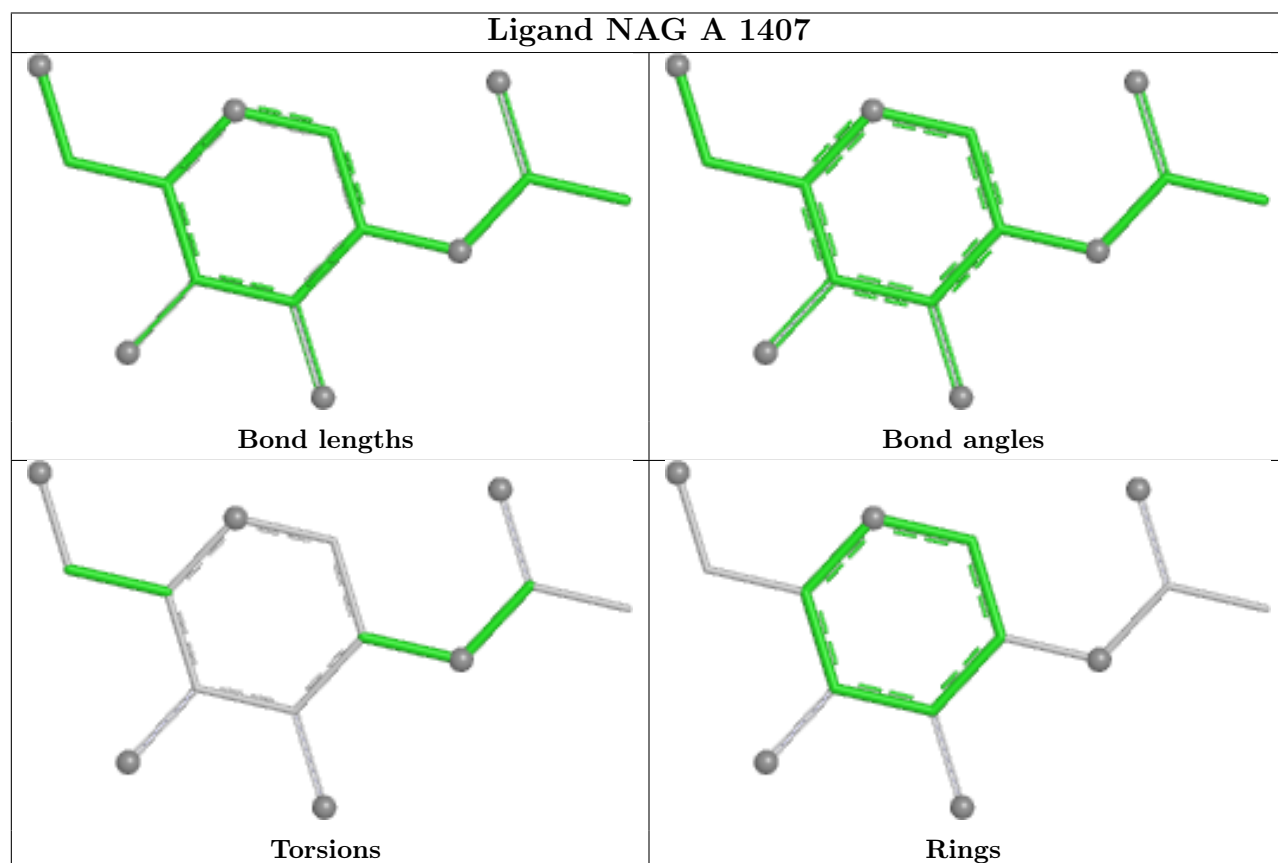
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



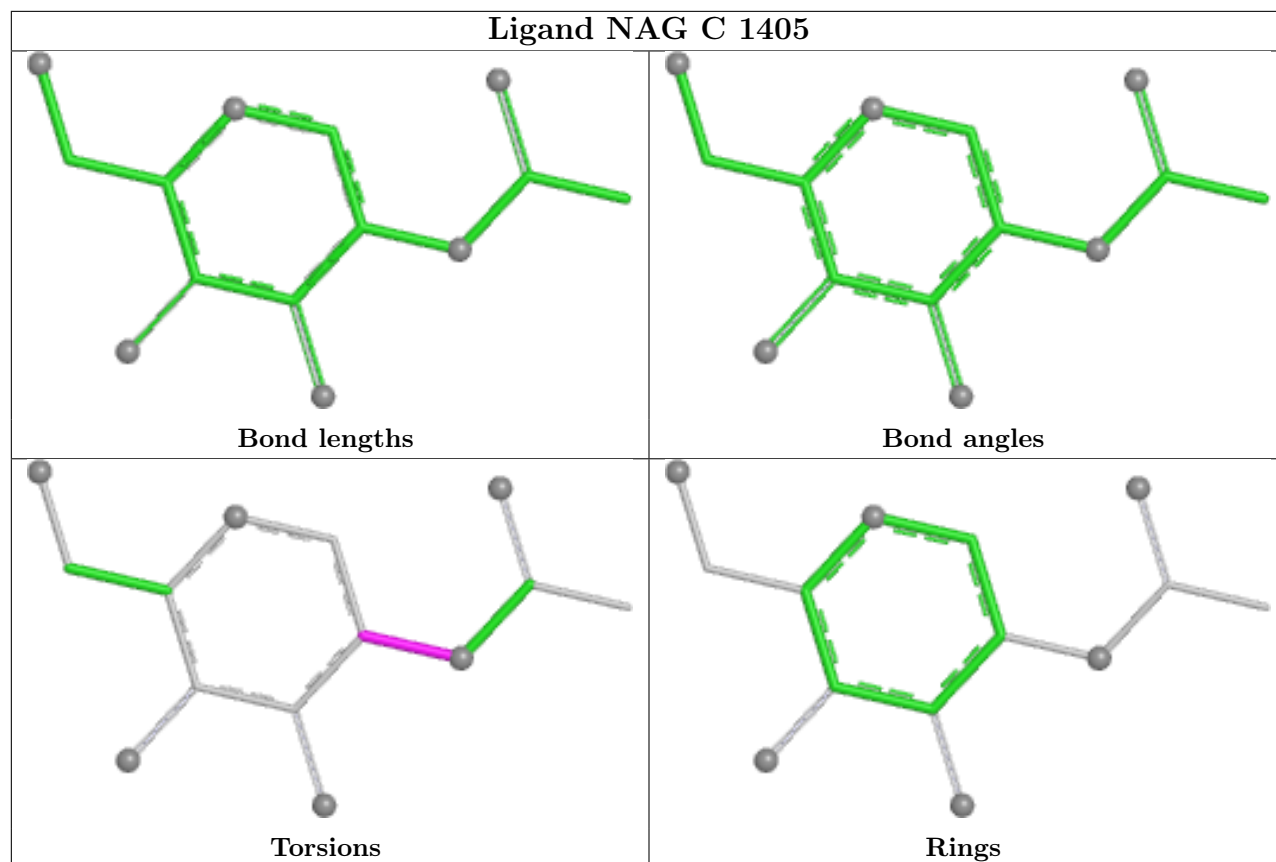
Ligand NAG B 1403



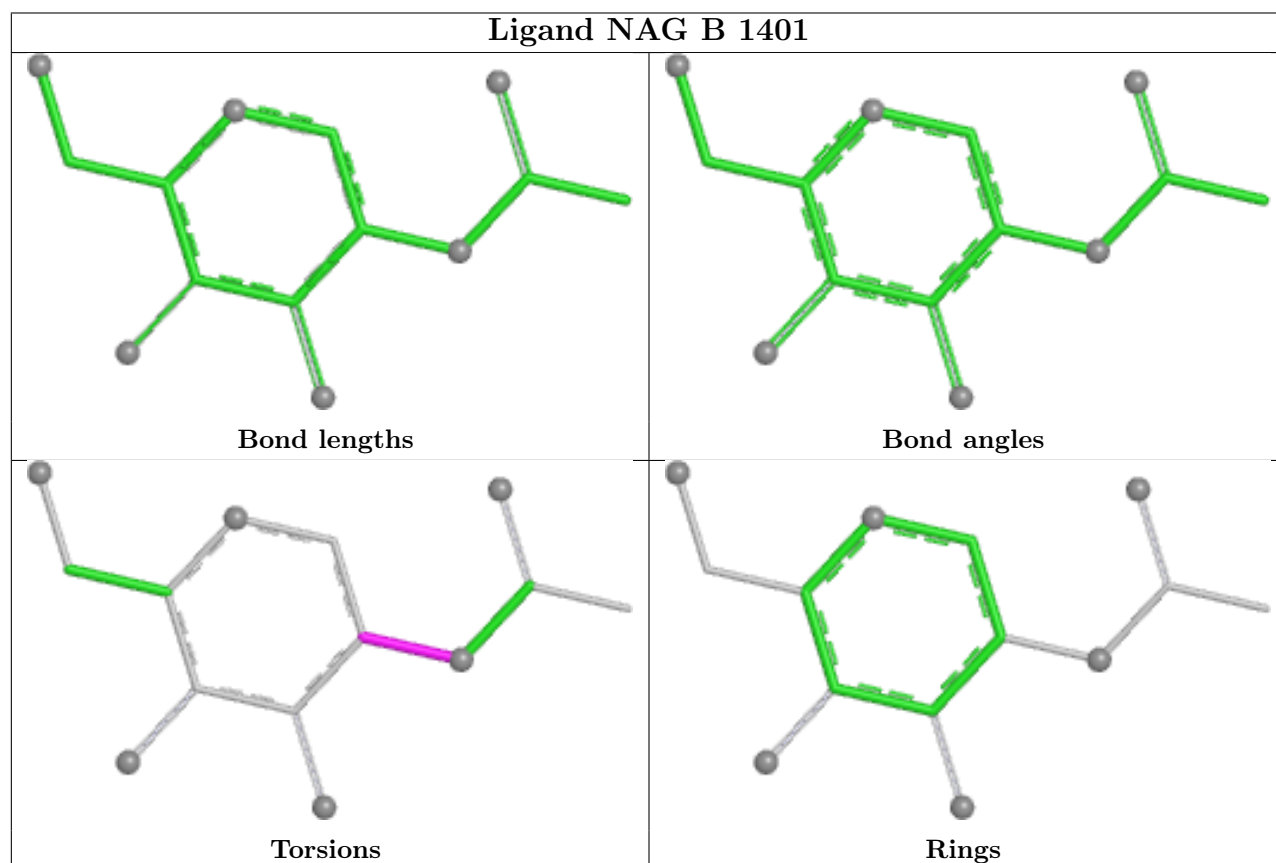
Ligand NAG A 1407



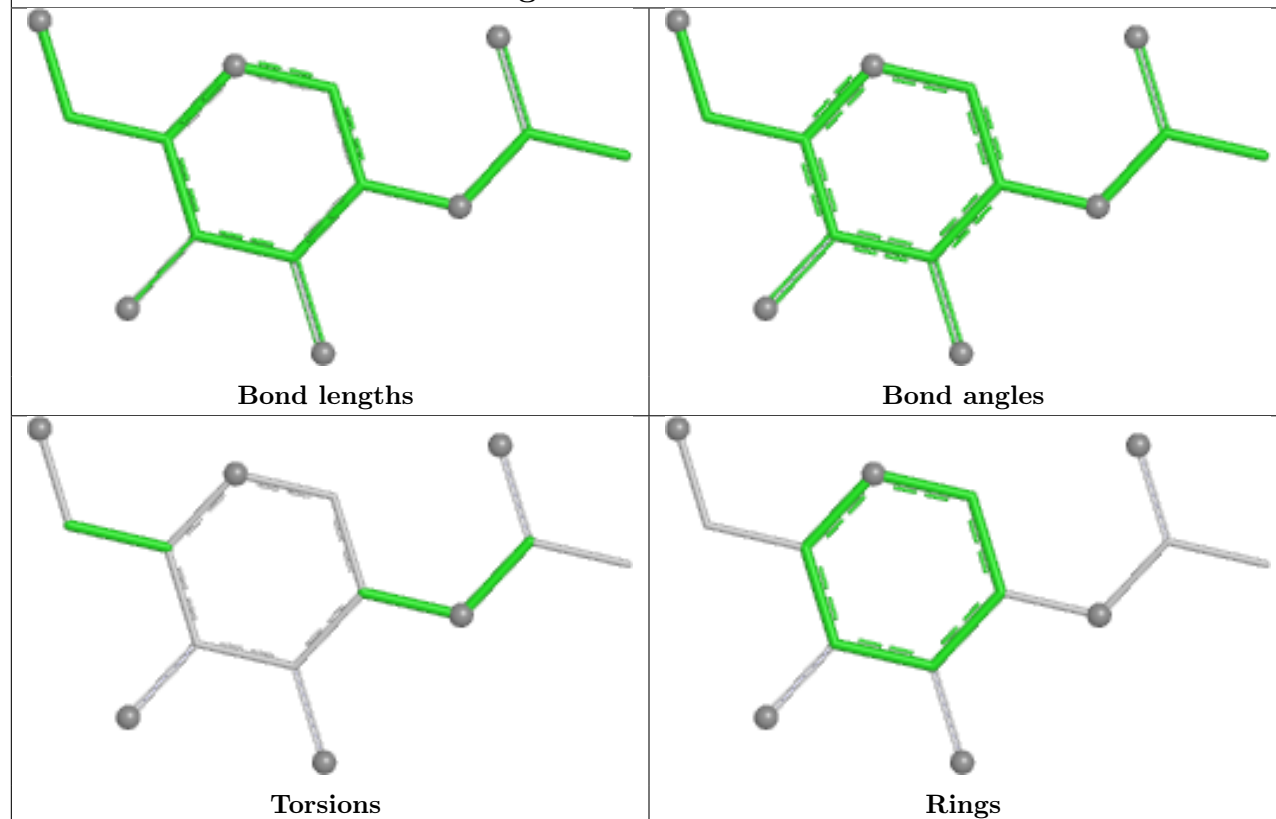
Ligand NAG C 1405



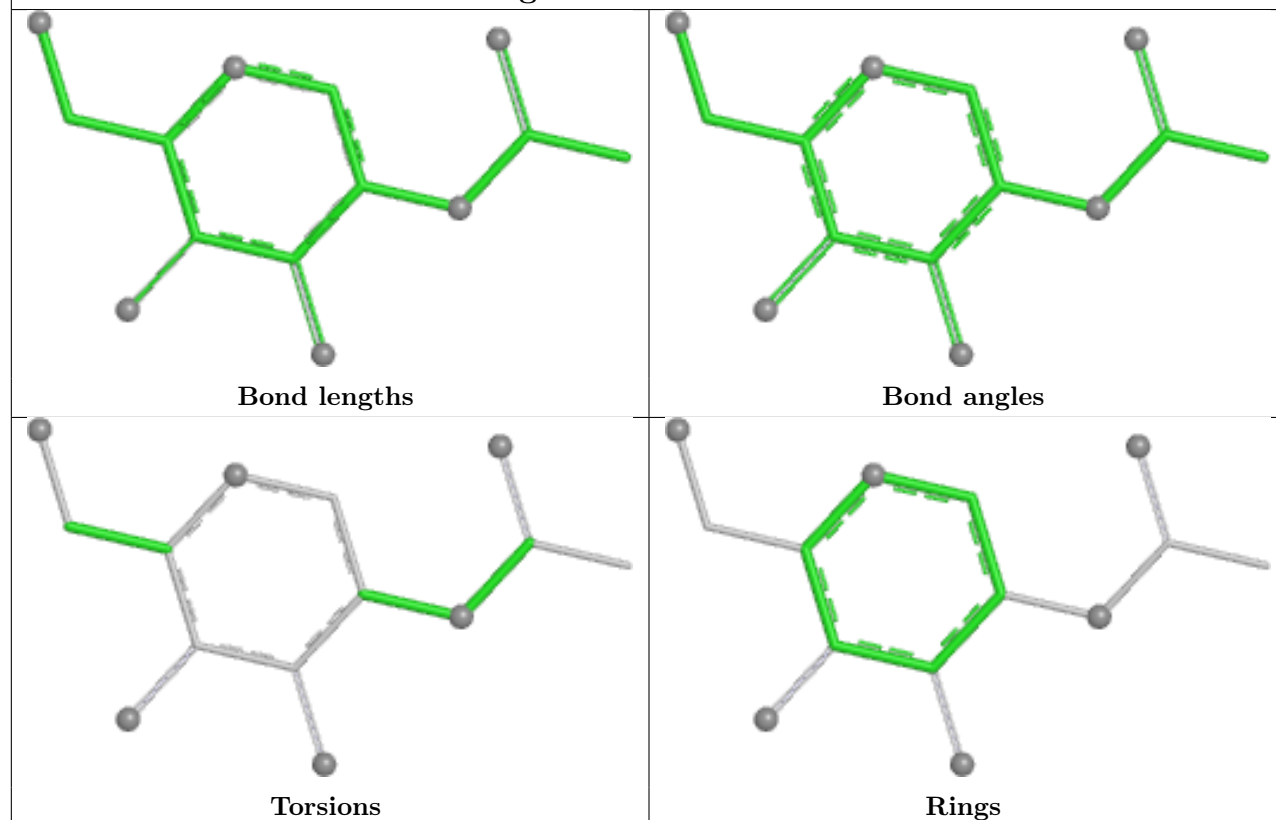
Ligand NAG B 1401



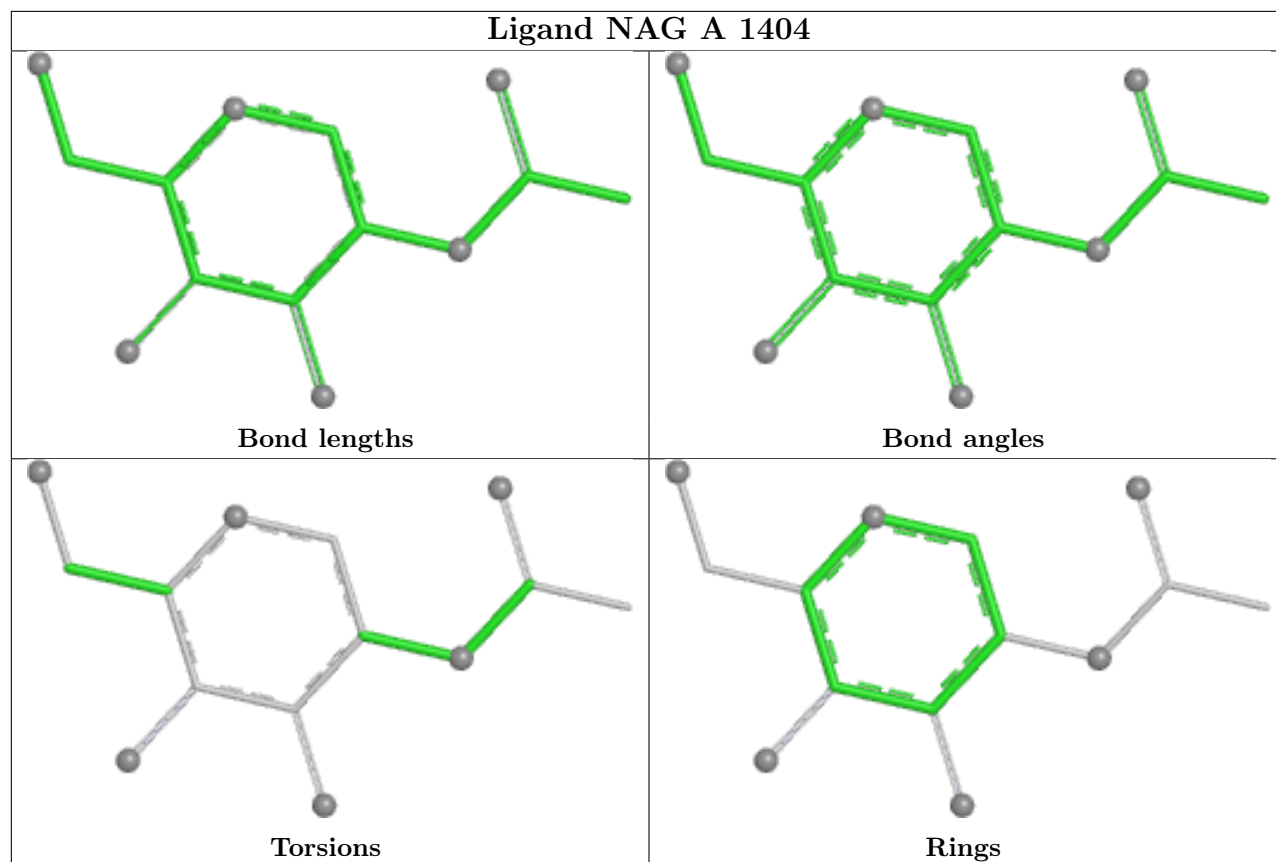
Ligand NAG A 1403



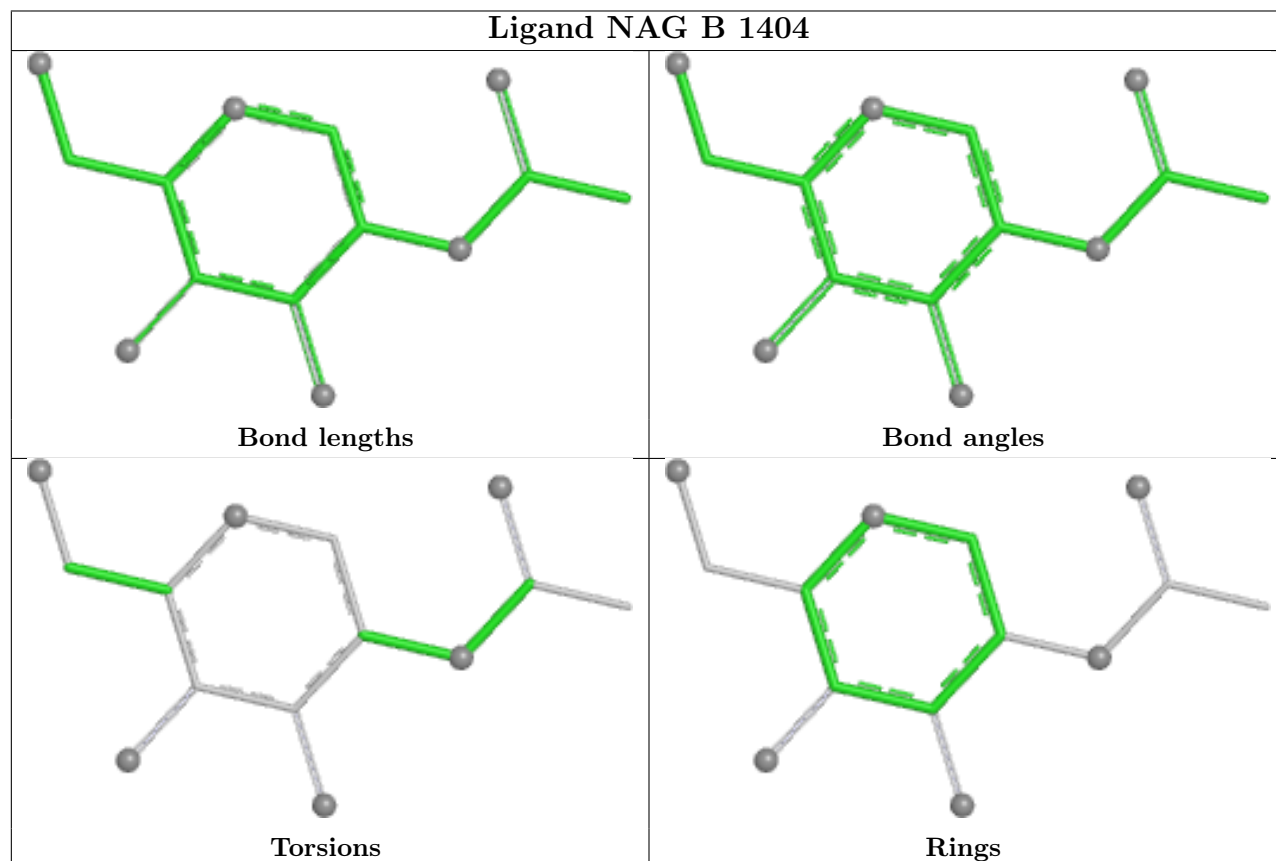
Ligand NAG C 1407



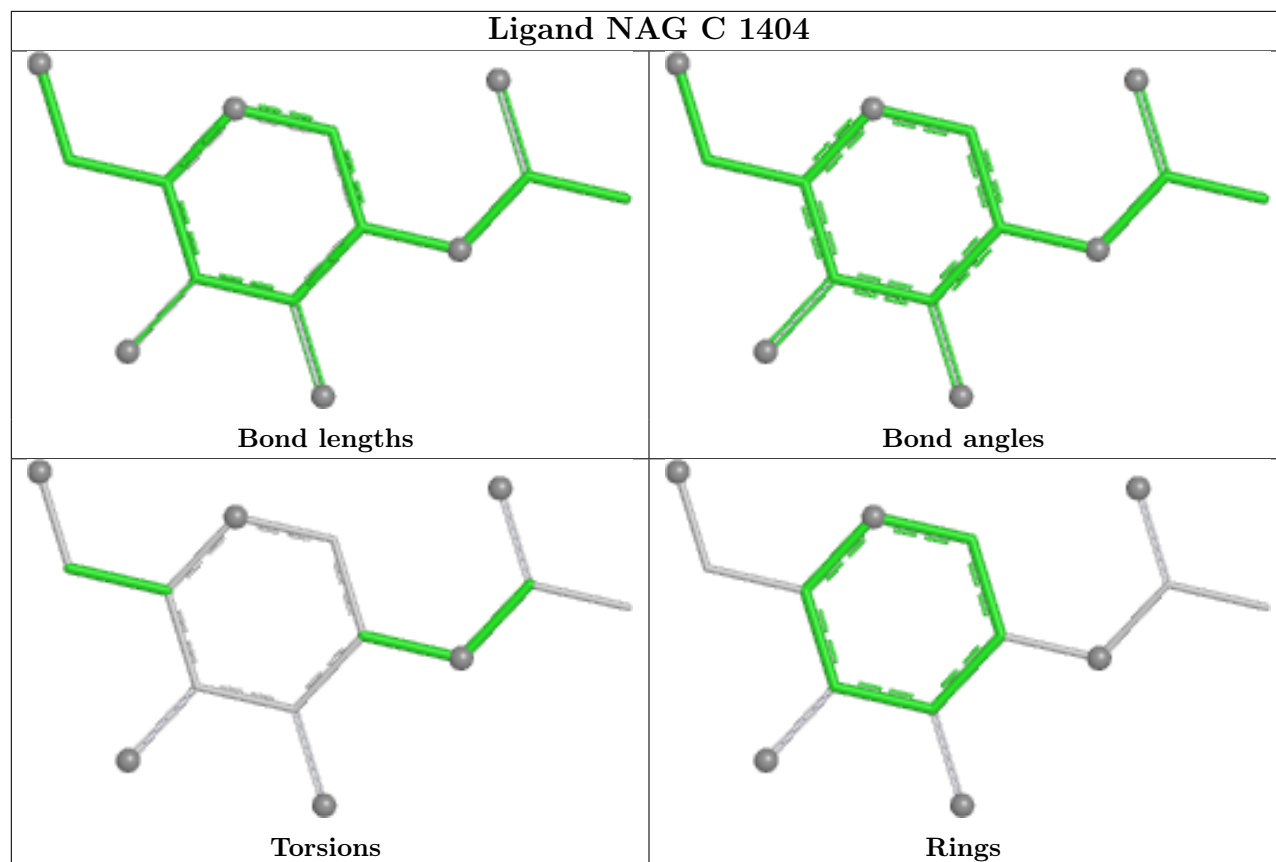
Ligand NAG A 1404



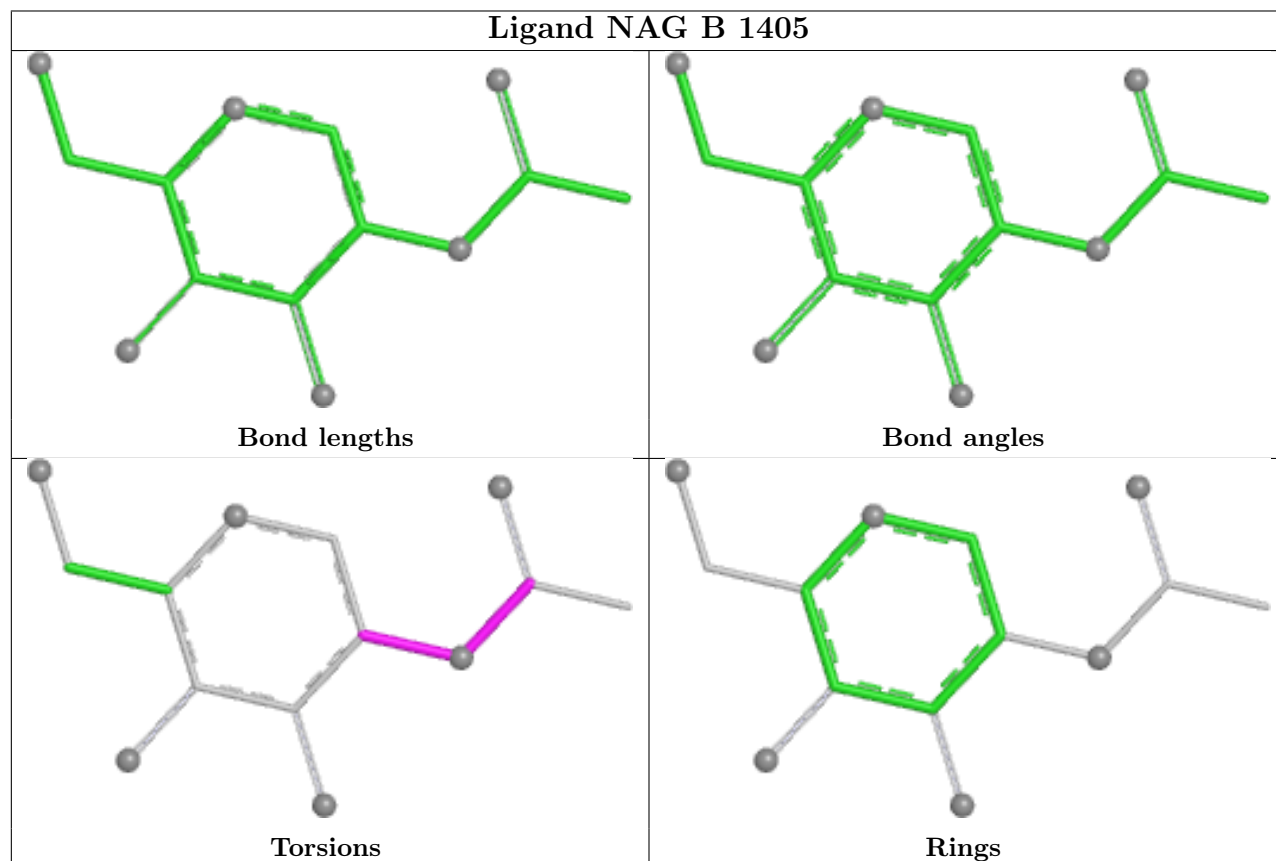
Ligand NAG B 1404

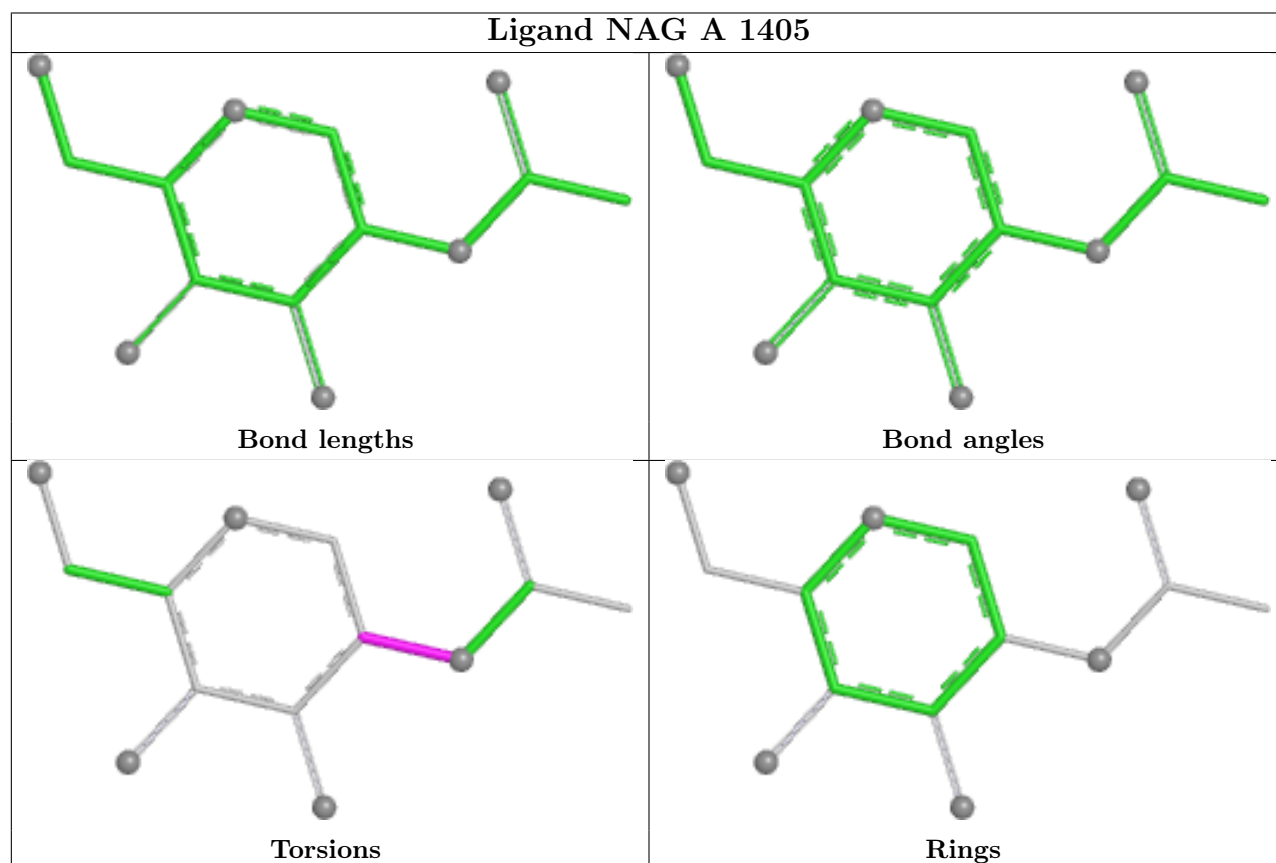
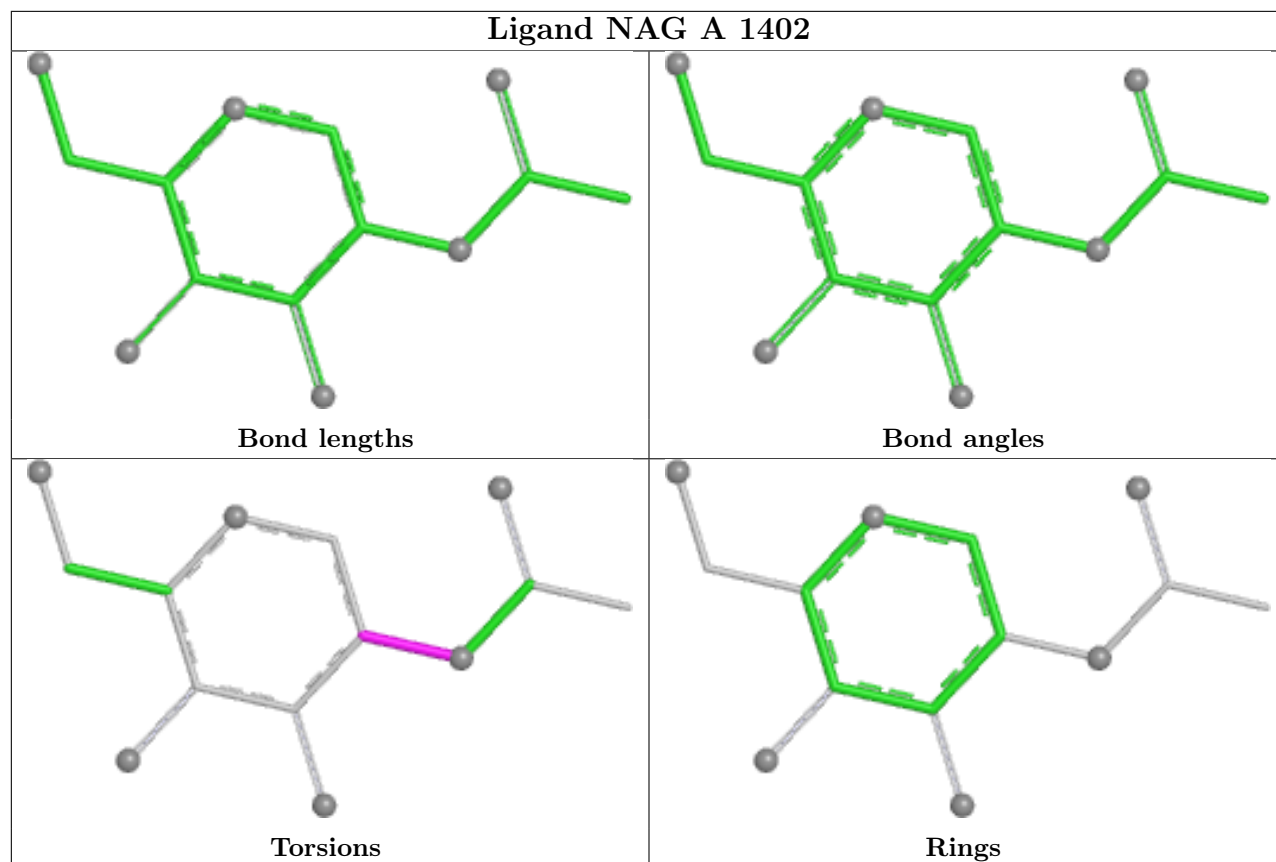


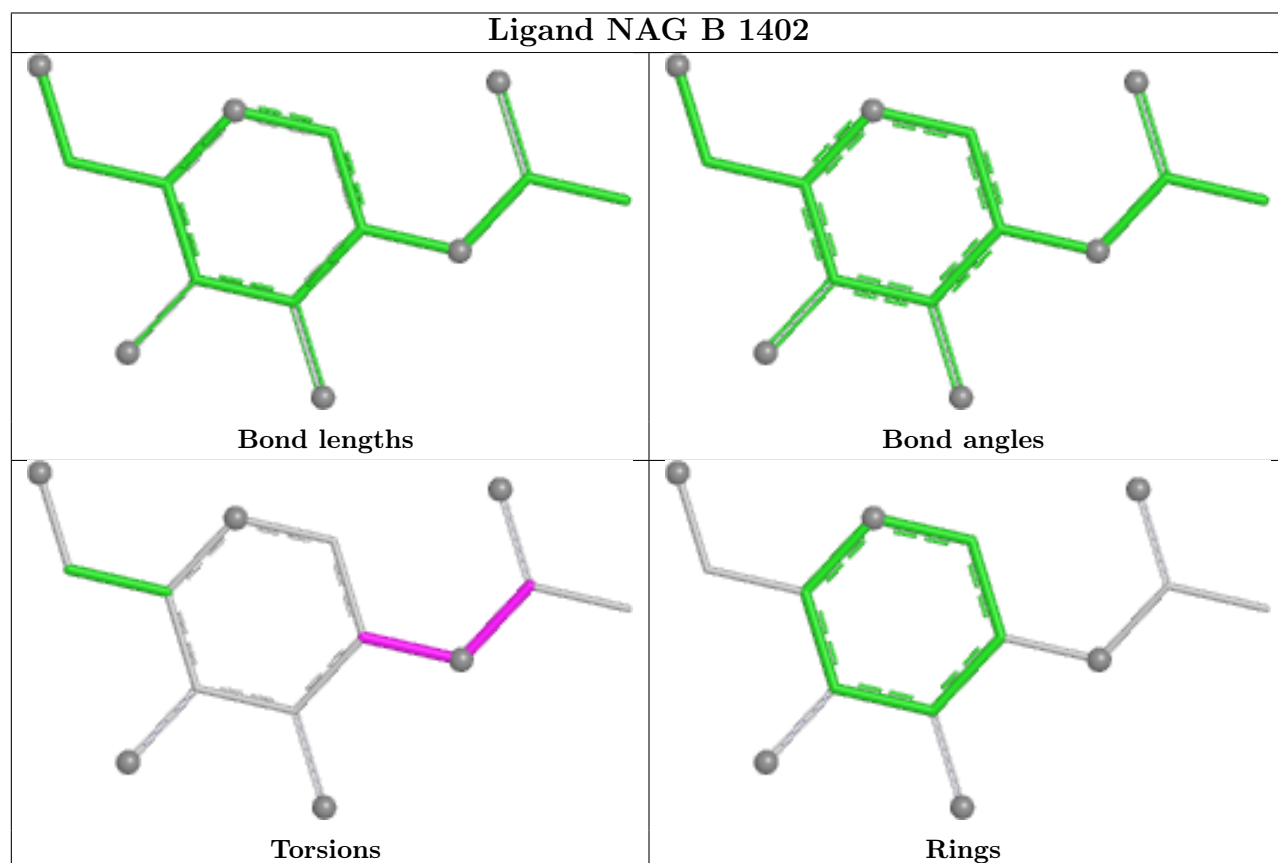
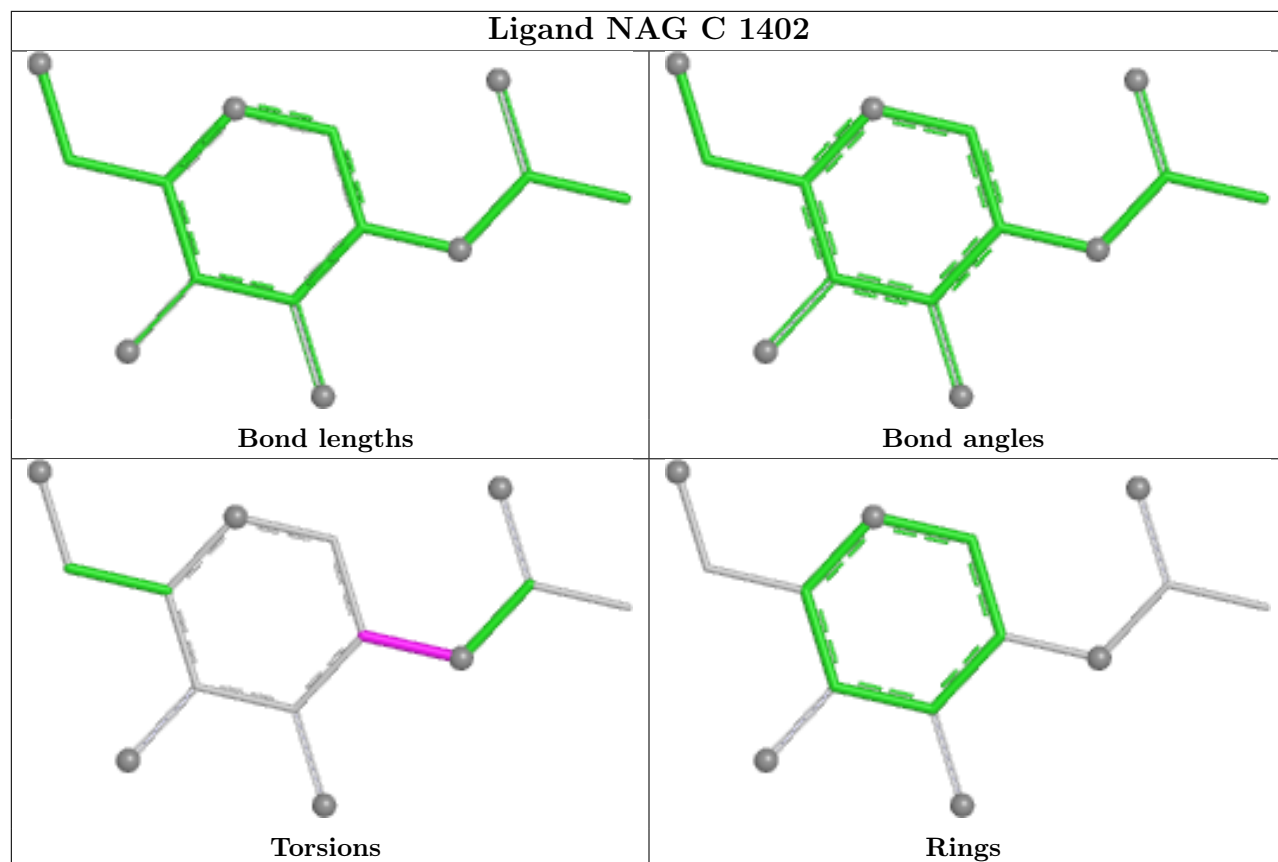
Ligand NAG C 1404

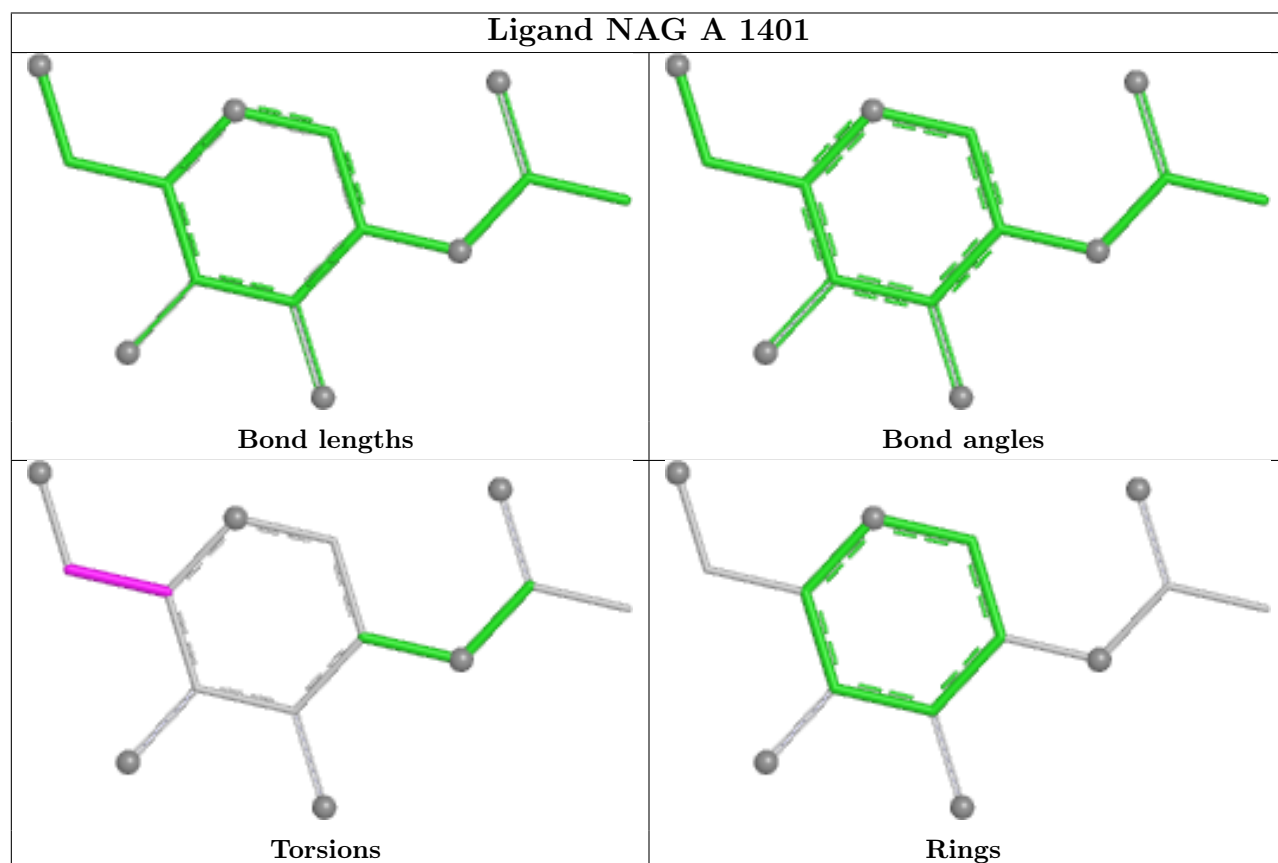
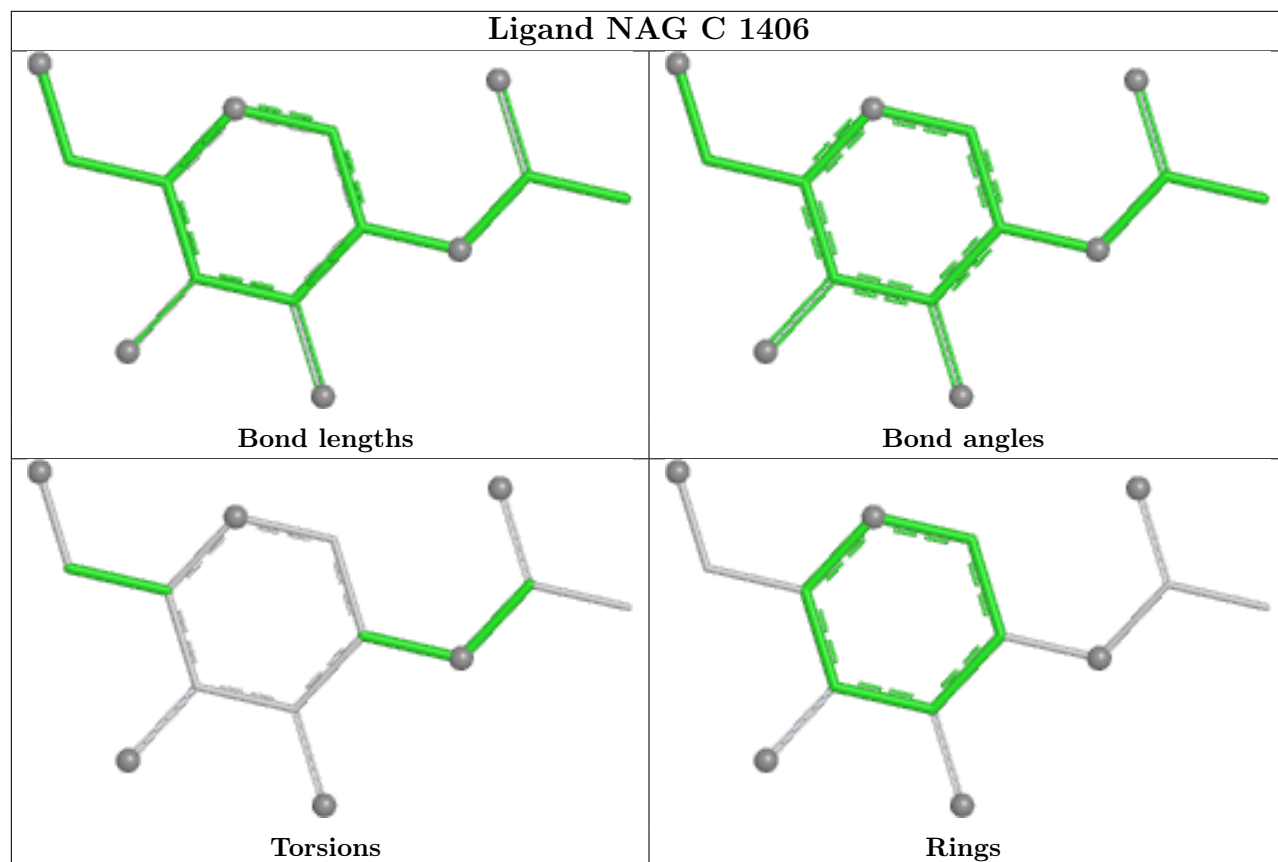


Ligand NAG B 1405

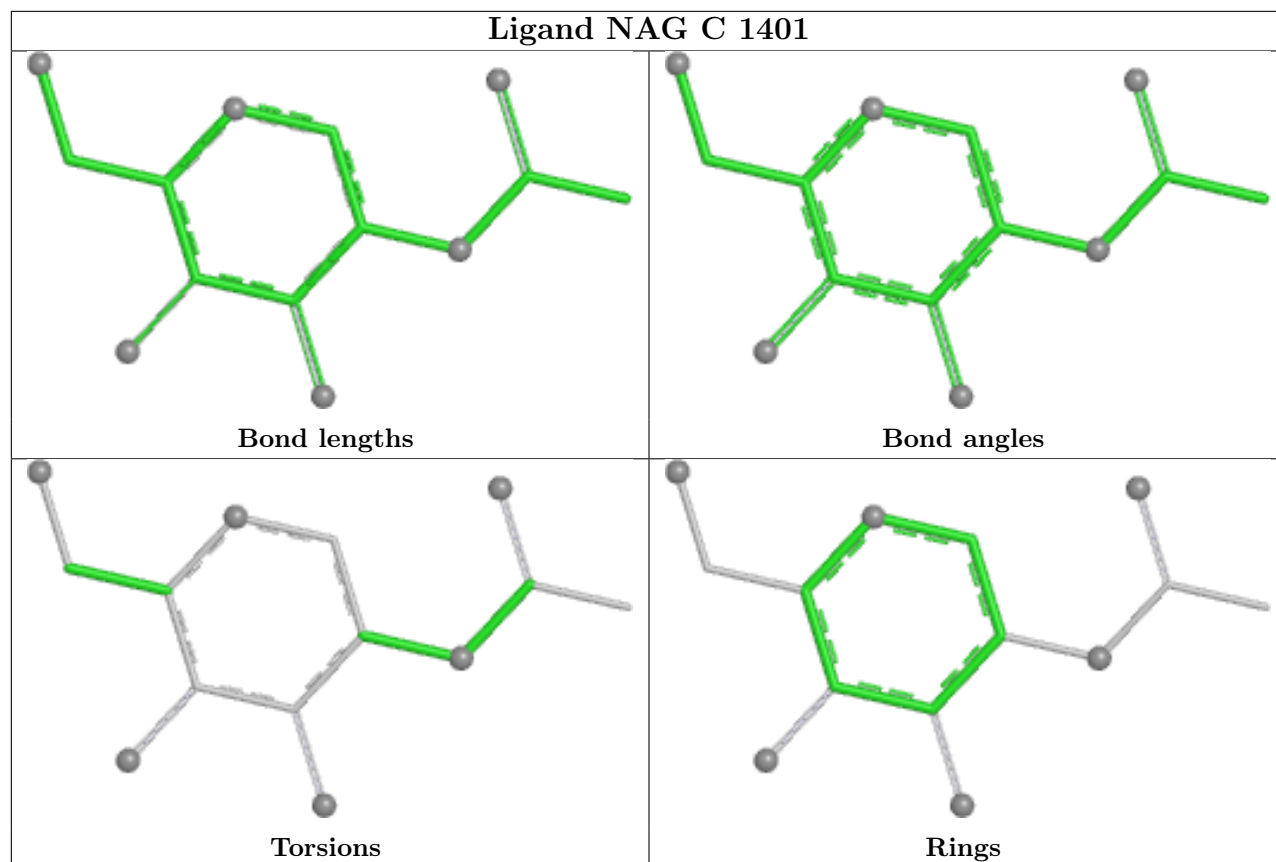




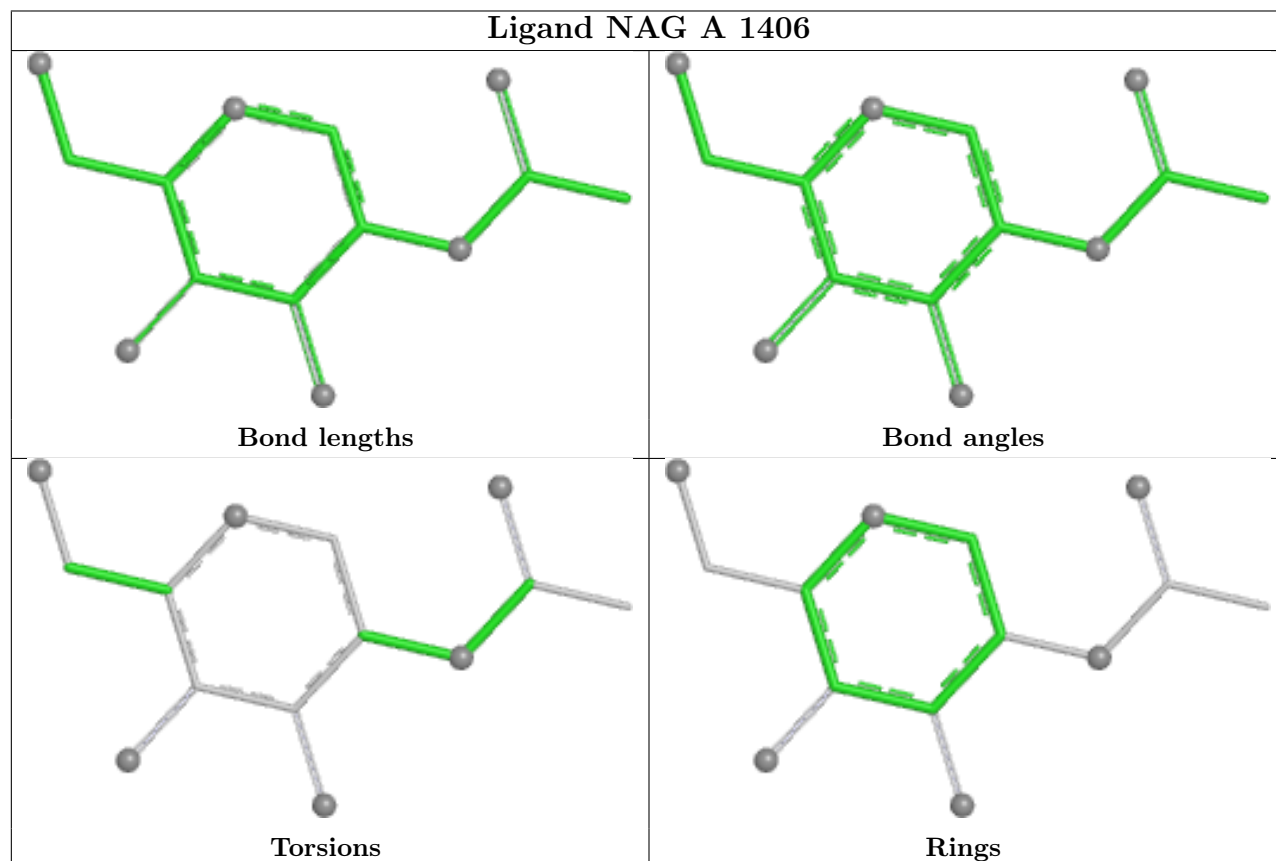




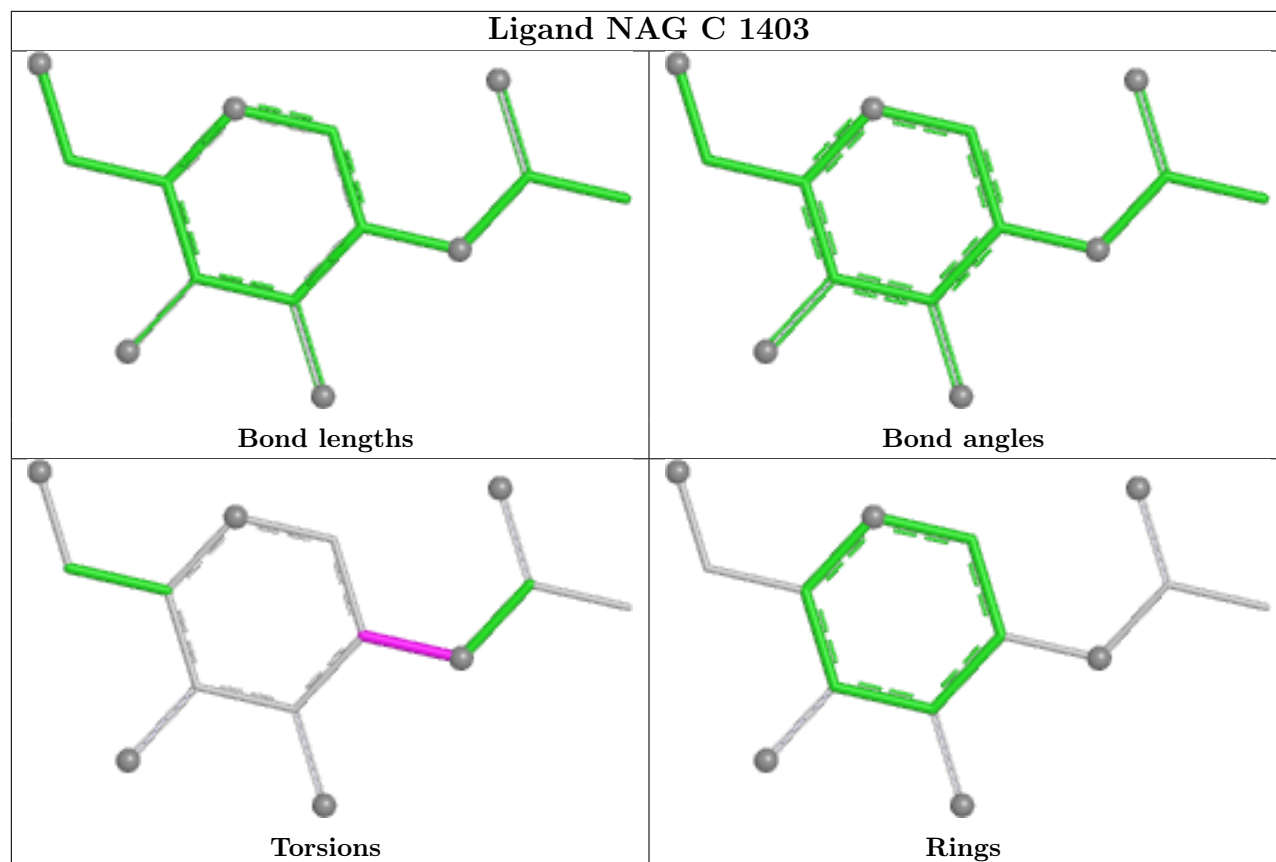
Ligand NAG C 1401



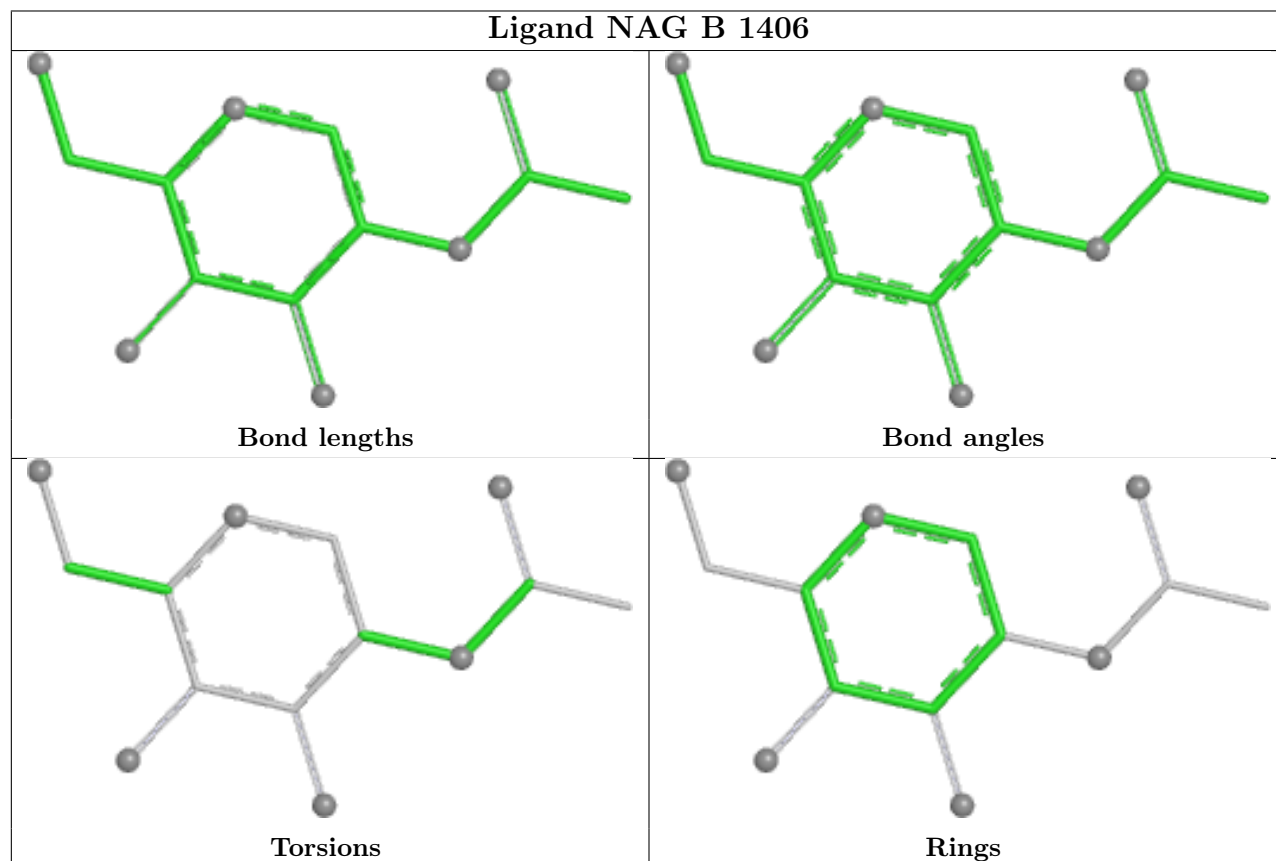
Ligand NAG A 1406



Ligand NAG C 1403



Ligand NAG B 1406



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.

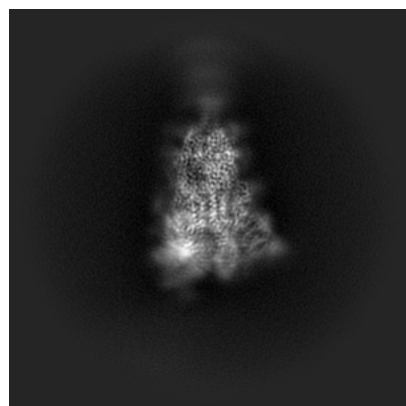
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-27206. These allow visual inspection of the internal detail of the map and identification of artifacts.

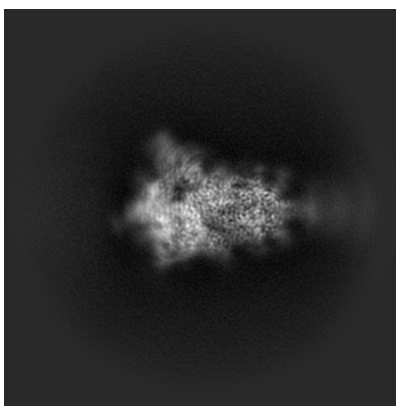
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

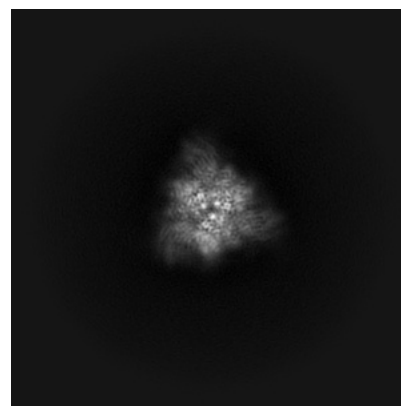
6.1.1 Primary map



X

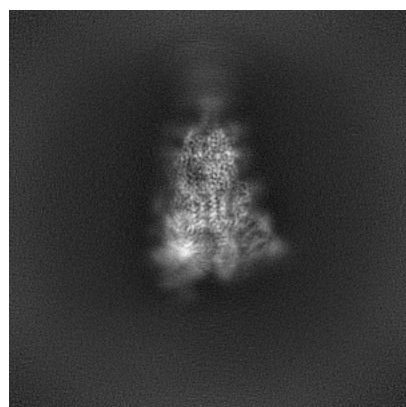


Y

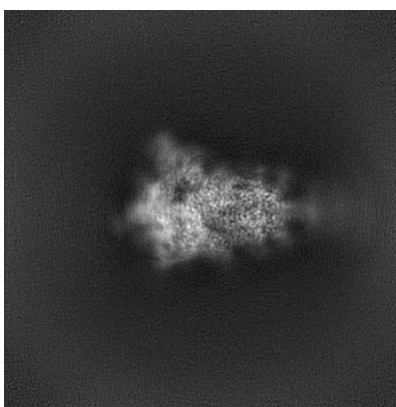


Z

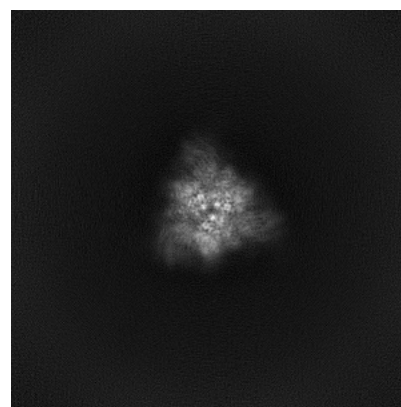
6.1.2 Raw map



X



Y

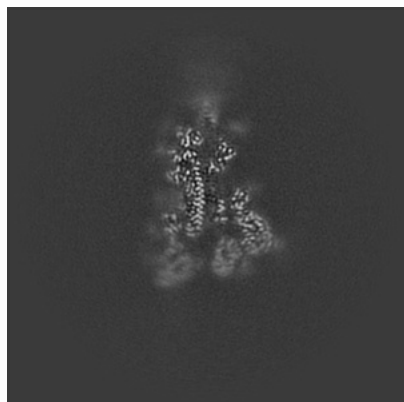


Z

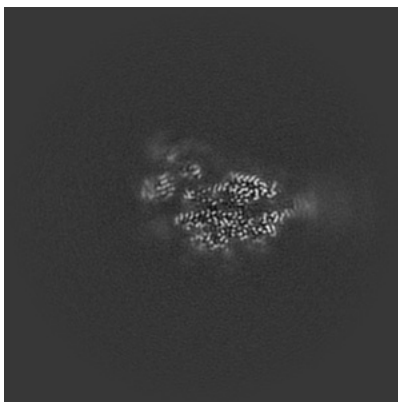
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

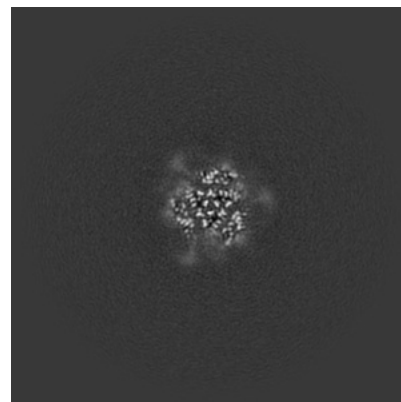
6.2.1 Primary map



X Index: 240

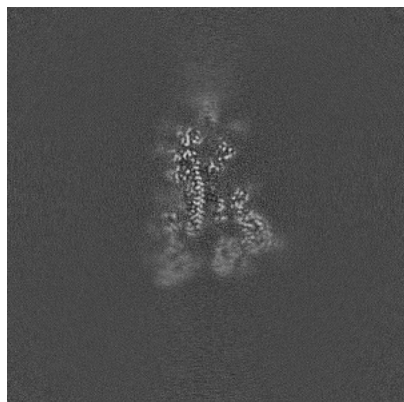


Y Index: 240

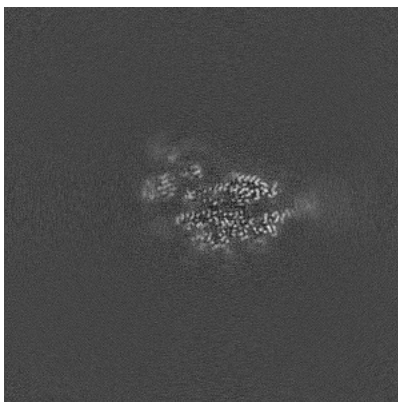


Z Index: 240

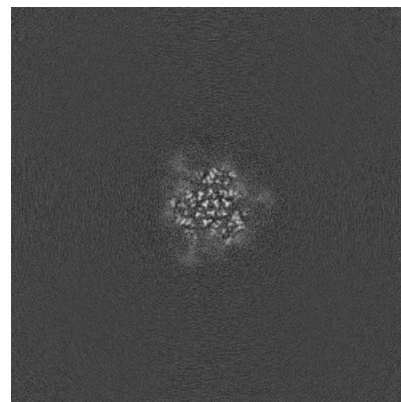
6.2.2 Raw map



X Index: 240



Y Index: 240

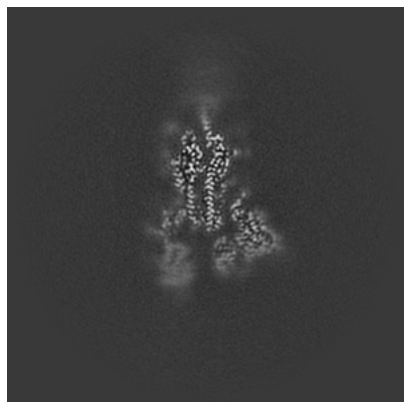


Z Index: 240

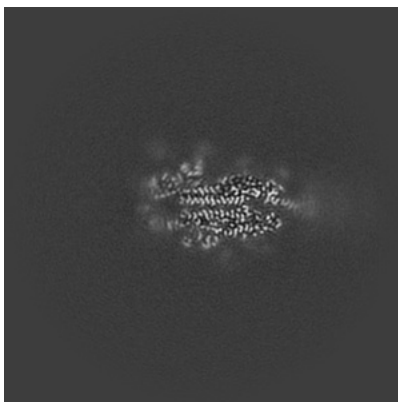
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

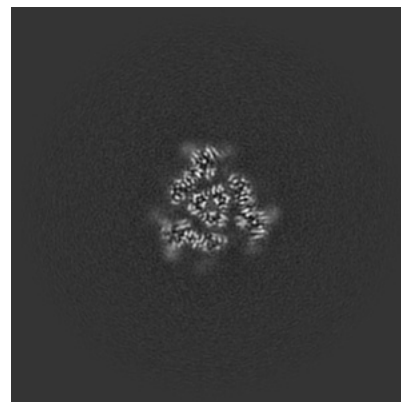
6.3.1 Primary map



X Index: 231

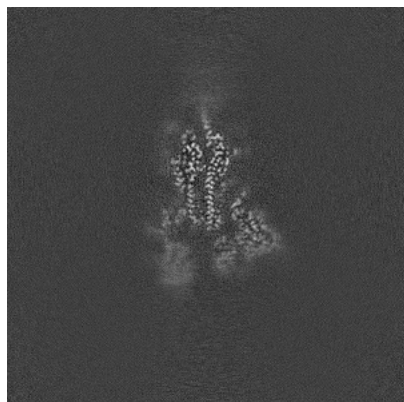


Y Index: 246



Z Index: 220

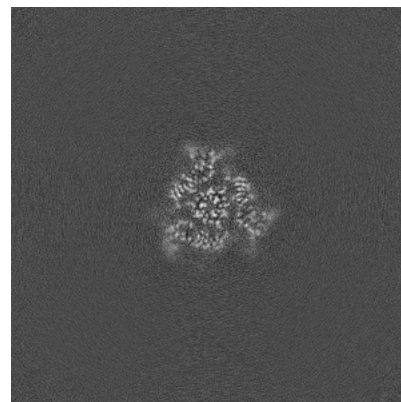
6.3.2 Raw map



X Index: 231



Y Index: 246

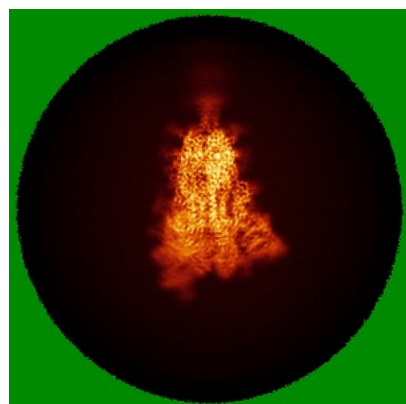


Z Index: 226

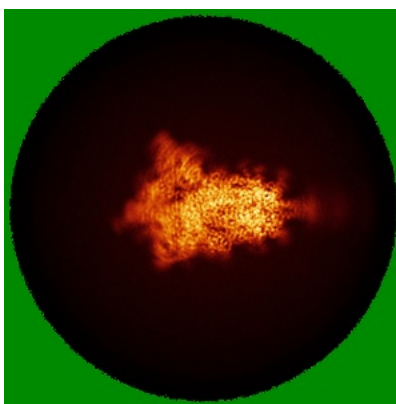
The images above show the largest variance slices of the map in three orthogonal directions.

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

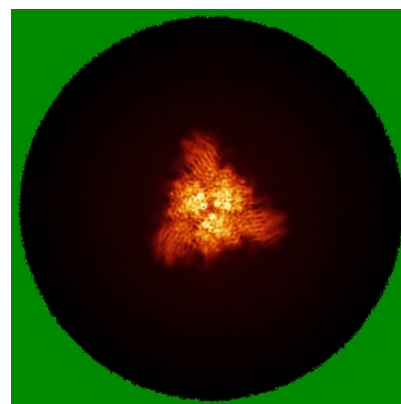
6.4.1 Primary map



X

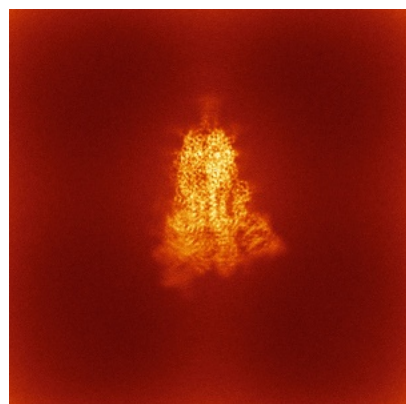


Y

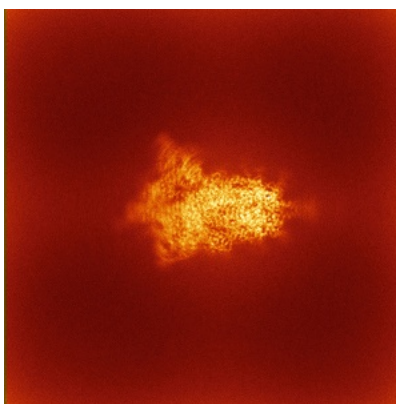


Z

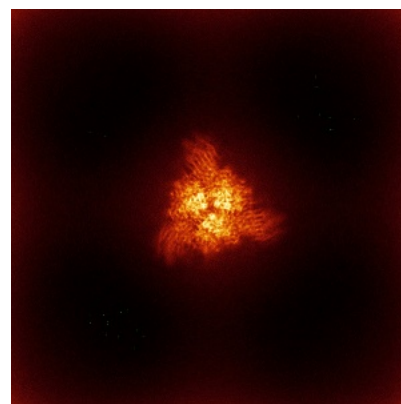
6.4.2 Raw map



X



Y

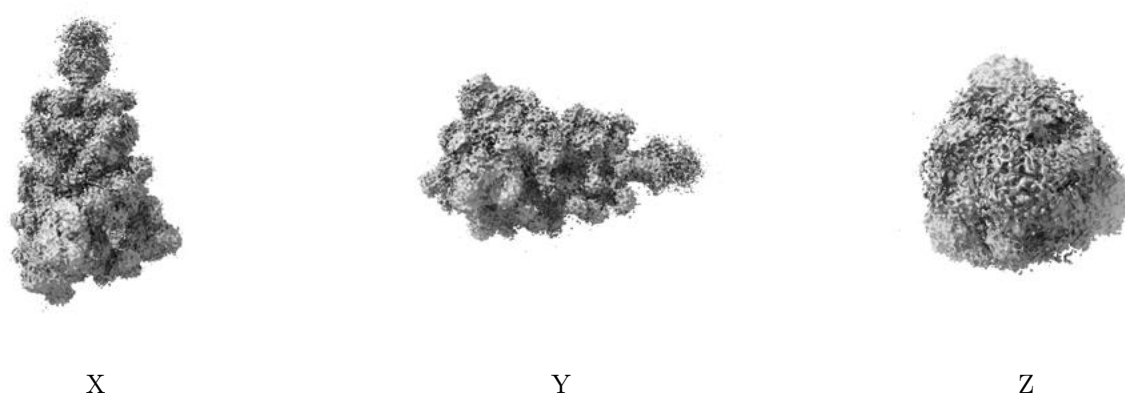


Z

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.

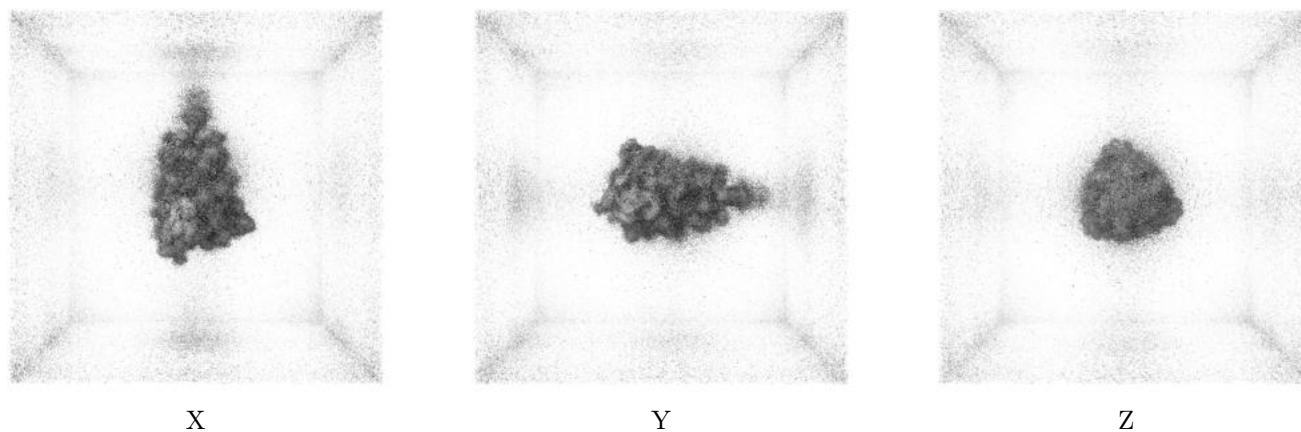
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.03. 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.

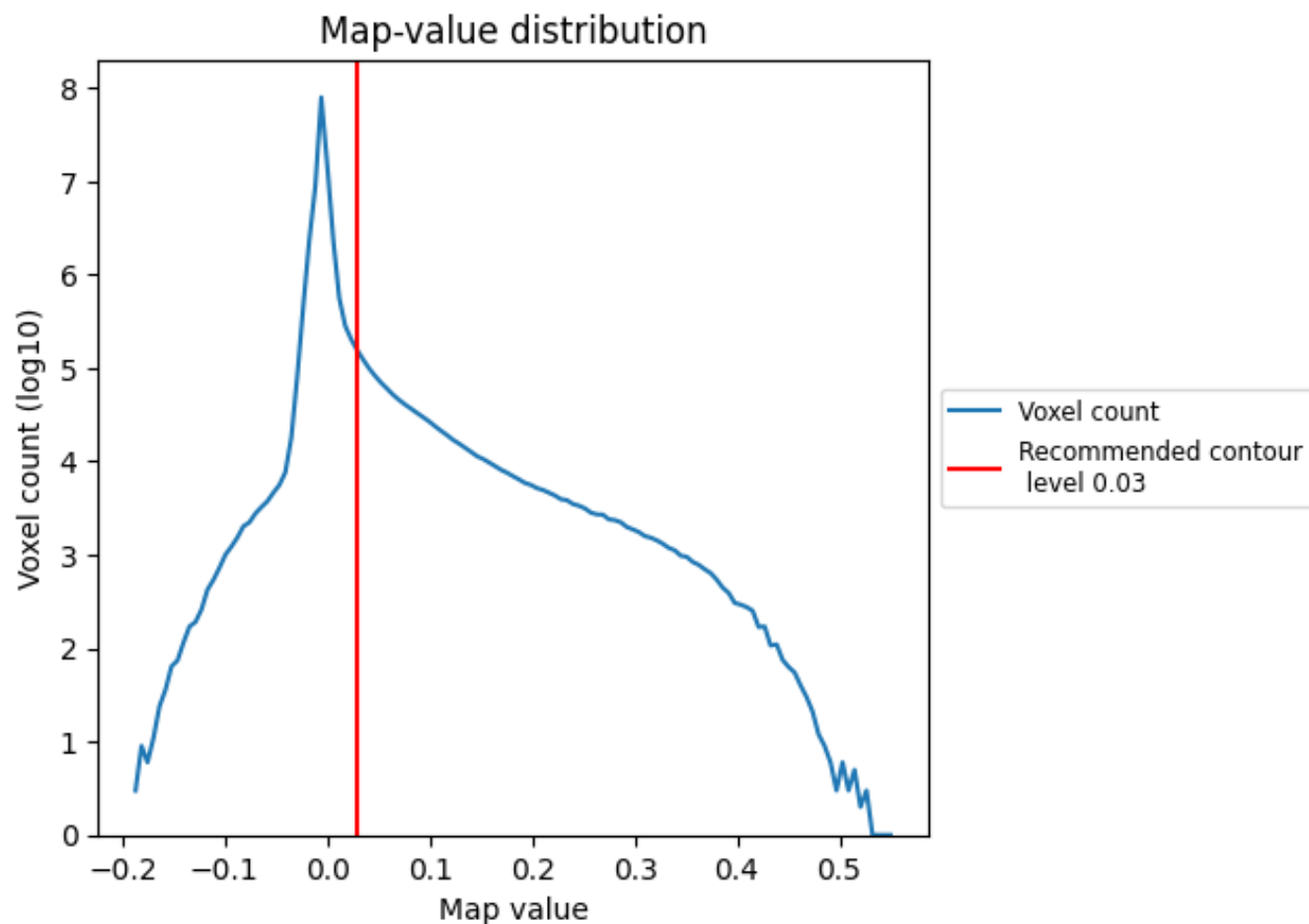
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

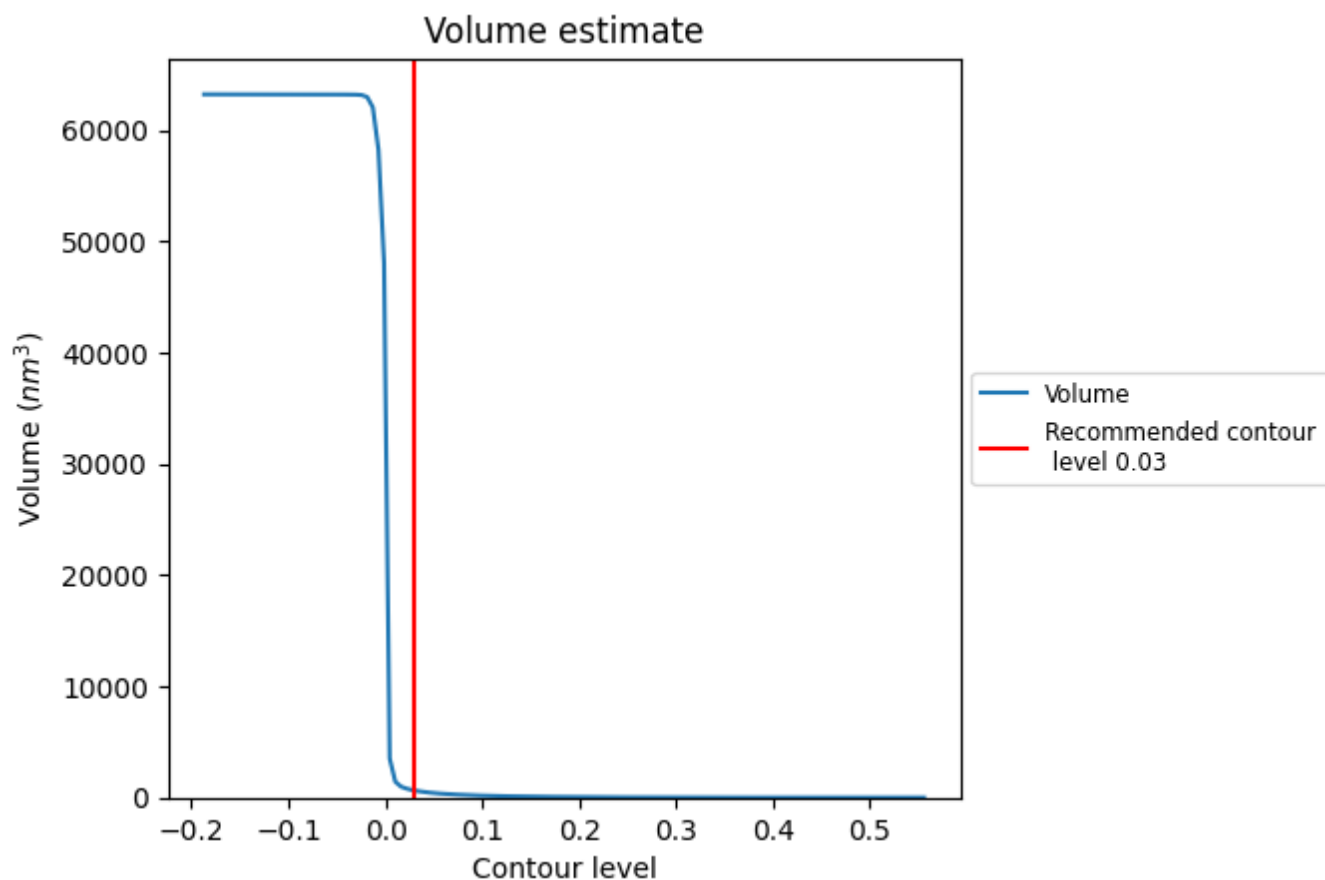
This section contains the results of statistical analysis of the map.

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.

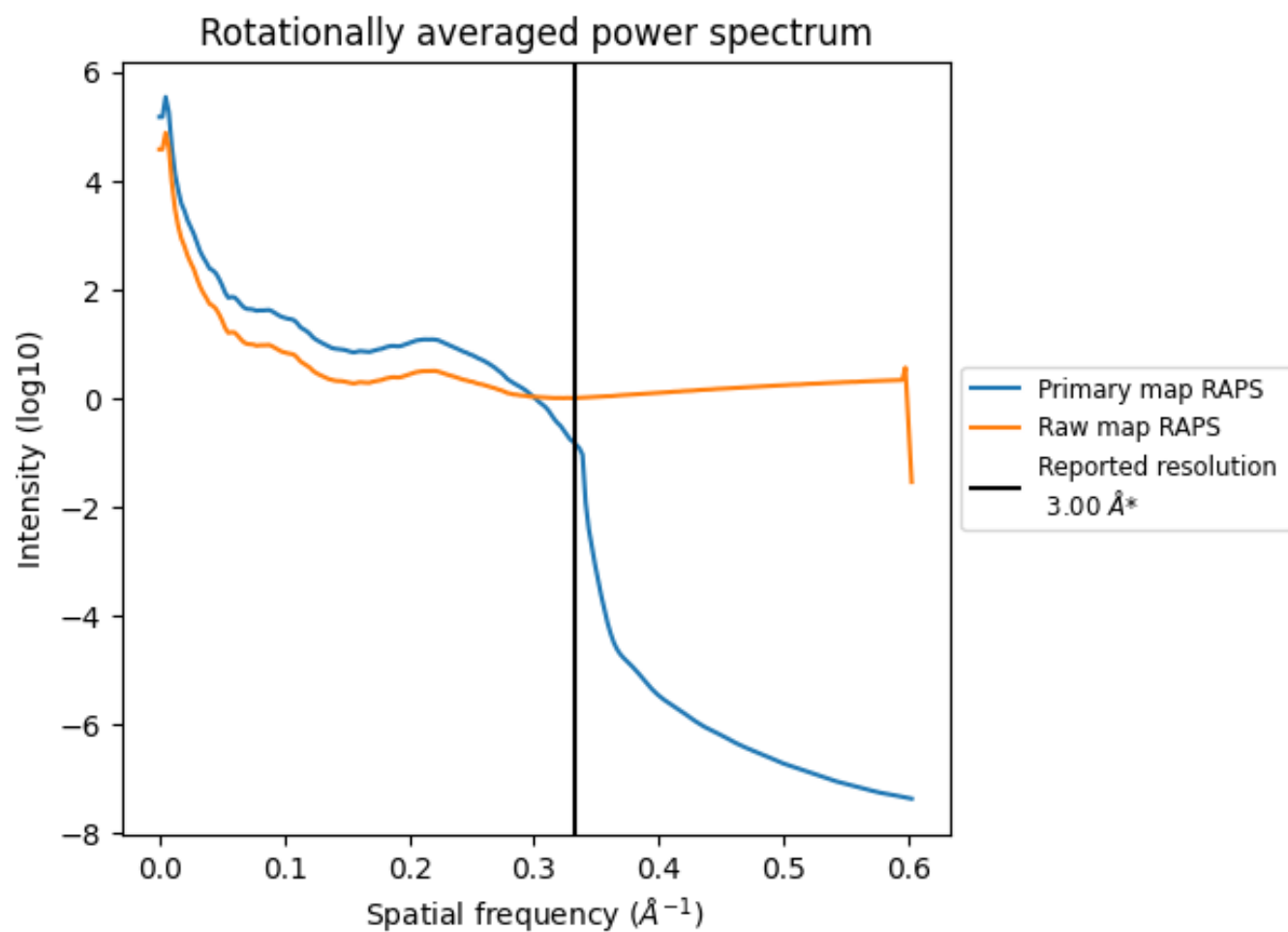
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 640 nm³; this corresponds to an approximate mass of 578 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 ⓘ

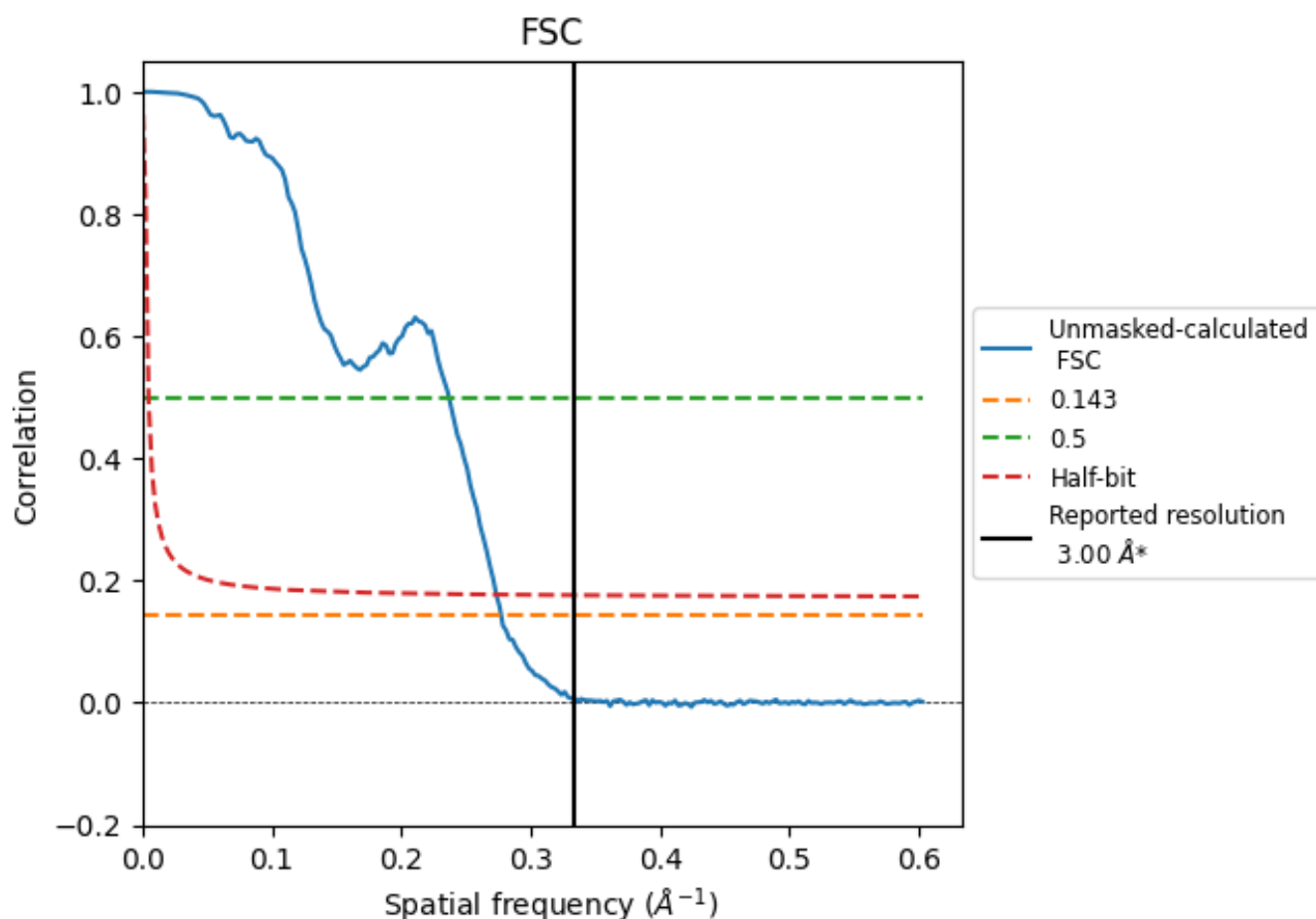


*Reported resolution corresponds to spatial frequency of 0.333 $\text{\AA}^{-1}$

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 0.333 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

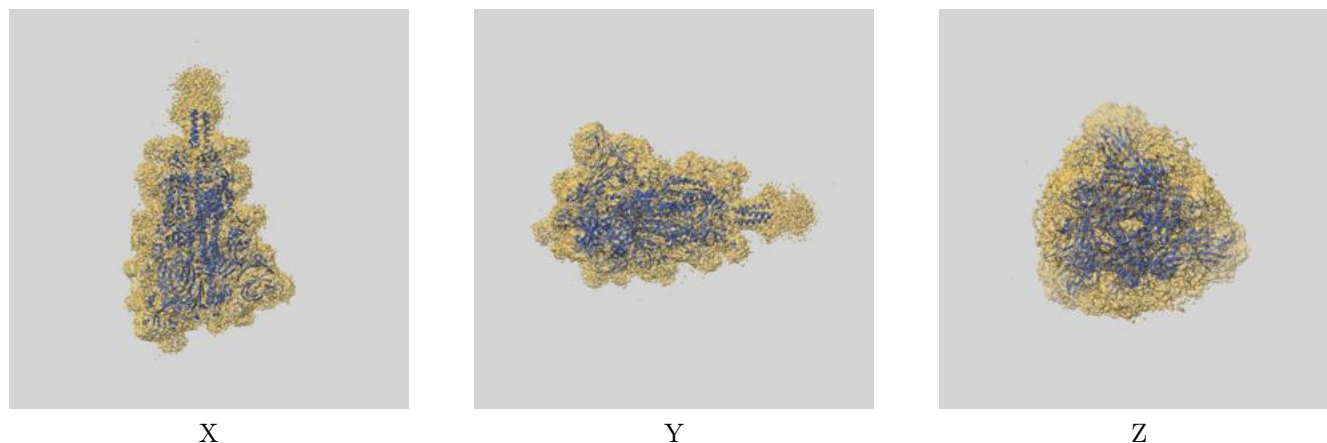
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.00	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.61	4.22	3.66

*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 3.61 differs from the reported value 3.0 by more than 10 %

9 Map-model fit [i](#)

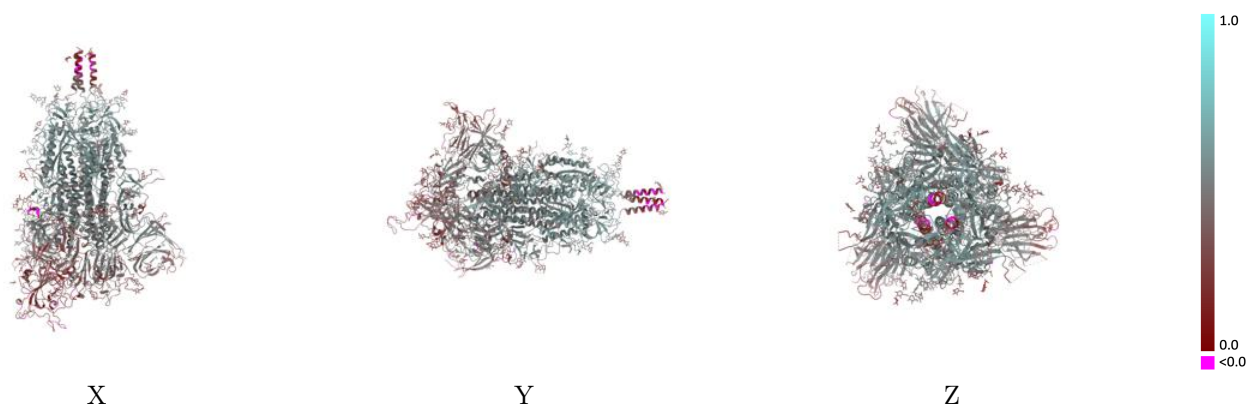
This section contains information regarding the fit between EMDB map EMD-27206 and PDB model 8D56. Per-residue inclusion information can be found in [section 3](#) on [page 14](#).

9.1 Map-model overlay [i](#)



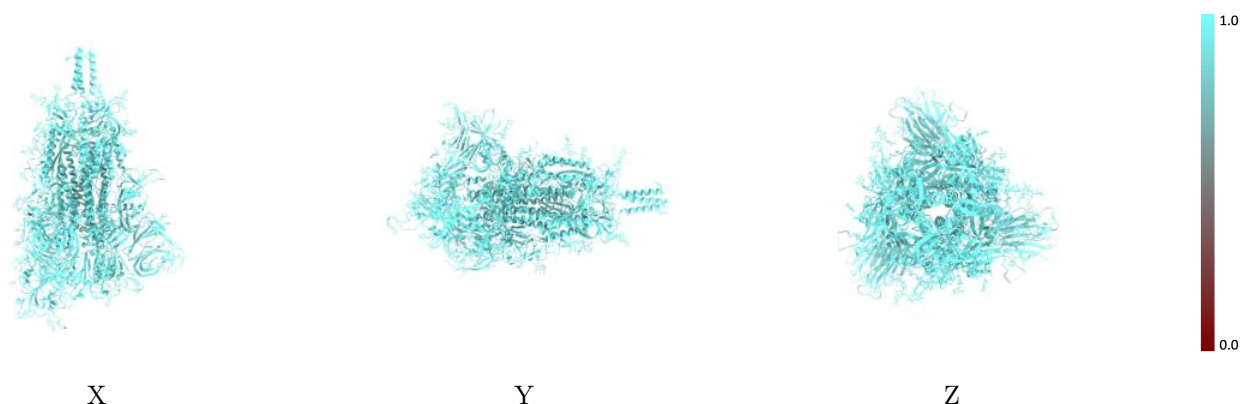
The images above show the 3D surface view of the map at the recommended contour level 0.03 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



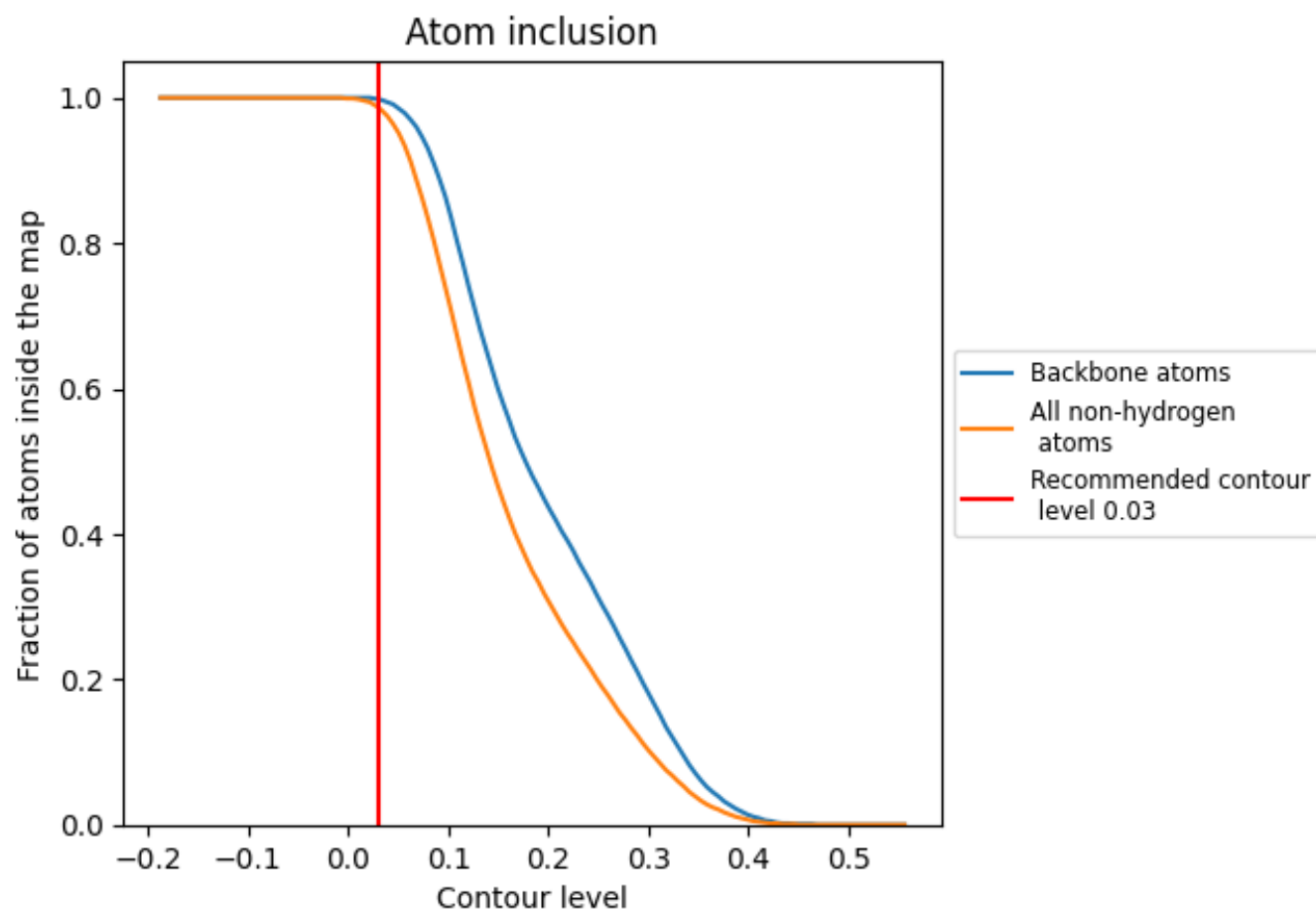
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.03).

9.4 Atom inclusion [i](#)



At the recommended contour level, 100% of all backbone atoms, 99% 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 (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9870	 0.4380
A	 0.9900	 0.4520
B	 0.9880	 0.4610
C	 0.9860	 0.4120
D	 1.0000	 0.2470
E	 1.0000	 0.2780
F	 0.6790	 0.1750
G	 1.0000	 0.2240
H	 0.9640	 0.3030
I	 1.0000	 0.5100
J	 1.0000	 0.4530
K	 1.0000	 0.4380
L	 1.0000	 0.4360
M	 1.0000	 0.4000
N	 1.0000	 0.2840
O	 1.0000	 0.2920
P	 1.0000	 0.3150
Q	 1.0000	 0.3370
R	 1.0000	 0.2160
S	 1.0000	 0.4930
T	 1.0000	 0.4740
U	 0.9740	 0.4490
V	 0.9740	 0.4160
W	 1.0000	 0.3600
X	 0.9740	 0.2000
Y	 0.8930	 0.1650
Z	 0.9640	 0.2410
a	 1.0000	 0.2600
b	 1.0000	 0.2510
c	 1.0000	 0.4720
d	 1.0000	 0.4310
e	 0.9470	 0.4140
f	 1.0000	 0.4350
g	 1.0000	 0.4480

