



Full wwPDB EM Validation Report ⓘ

Jun 10, 2025 – 12:20 PM JST

PDB ID : 9JMP / pdb_00009jmp
EMDB ID : EMD-61609
Title : Cryo-EM structure of GD-BatCoV (BtCoV/Ii/GD/2014-422) S-trimer
Authors : Yuan, H.; Xiong, X.; Gao, X.; Li, Z.; Wang, J.
Deposited on : 2024-09-20
Resolution : 3.10 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev118
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0rc1
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

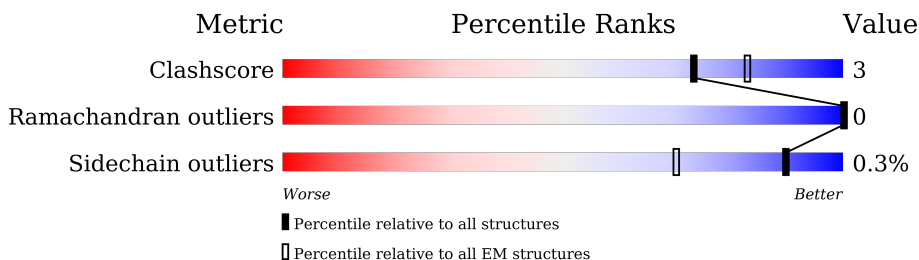
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.10 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	1373	<div> <div>72%</div> <div>6%</div> <div>22%</div> </div>
1	B	1373	<div> <div>72%</div> <div>6%</div> <div>22%</div> </div>
1	C	1373	<div> <div>71%</div> <div>7%</div> <div>22%</div> </div>
2	D	2	<div> <div>100%</div> </div>
2	F	2	<div> <div>100%</div> </div>
2	G	2	<div> <div>100%</div> </div>
2	I	2	<div> <div>100%</div> </div>
2	J	2	<div> <div>50%</div> <div>50%</div> </div>

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Mol	Chain	Length	Quality of chain
2	K	2	 100%
2	L	2	 100%
2	N	2	 50%50%
2	O	2	 100%
3	E	3	 100%
3	H	3	 100%
3	M	3	 67%33%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 25407 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,Fibritin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1065	Total	C	N	O	S	0	0
			8228	5218	1364	1596	50		
1	B	1065	Total	C	N	O	S	0	0
			8228	5218	1364	1596	50		
1	C	1065	Total	C	N	O	S	0	0
			8228	5218	1364	1596	50		

There are 177 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	885	PRO	ALA	engineered mutation	UNP A0A2R4KP93
A	962	PRO	SER	engineered mutation	UNP A0A2R4KP93
A	969	PRO	ALA	engineered mutation	UNP A0A2R4KP93
A	1012	PRO	ASN	engineered mutation	UNP A0A2R4KP93
A	1056	PRO	ALA	engineered mutation	UNP A0A2R4KP93
A	1057	PRO	VAL	engineered mutation	UNP A0A2R4KP93
A	1295	GLY	-	linker	UNP A0A2R4KP93
A	1296	SER	-	linker	UNP A0A2R4KP93
A	1318	LEU	PHE	conflict	UNP P10104
A	1324	LEU	-	expression tag	UNP P10104
A	1325	GLU	-	expression tag	UNP P10104
A	1326	VAL	-	expression tag	UNP P10104
A	1327	LEU	-	expression tag	UNP P10104
A	1328	PHE	-	expression tag	UNP P10104
A	1329	GLN	-	expression tag	UNP P10104
A	1330	GLY	-	expression tag	UNP P10104
A	1331	PRO	-	expression tag	UNP P10104
A	1332	GLY	-	expression tag	UNP P10104
A	1333	HIS	-	expression tag	UNP P10104
A	1334	HIS	-	expression tag	UNP P10104
A	1335	HIS	-	expression tag	UNP P10104
A	1336	HIS	-	expression tag	UNP P10104
A	1337	HIS	-	expression tag	UNP P10104
A	1338	HIS	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1339	HIS	-	expression tag	UNP P10104
A	1340	HIS	-	expression tag	UNP P10104
A	1341	SER	-	expression tag	UNP P10104
A	1342	ALA	-	expression tag	UNP P10104
A	1343	TRP	-	expression tag	UNP P10104
A	1344	SER	-	expression tag	UNP P10104
A	1345	HIS	-	expression tag	UNP P10104
A	1346	PRO	-	expression tag	UNP P10104
A	1347	GLN	-	expression tag	UNP P10104
A	1348	PHE	-	expression tag	UNP P10104
A	1349	GLU	-	expression tag	UNP P10104
A	1350	LYS	-	expression tag	UNP P10104
A	1351	GLY	-	expression tag	UNP P10104
A	1352	GLY	-	expression tag	UNP P10104
A	1353	GLY	-	expression tag	UNP P10104
A	1354	SER	-	expression tag	UNP P10104
A	1355	GLY	-	expression tag	UNP P10104
A	1356	GLY	-	expression tag	UNP P10104
A	1357	GLY	-	expression tag	UNP P10104
A	1358	GLY	-	expression tag	UNP P10104
A	1359	SER	-	expression tag	UNP P10104
A	1360	GLY	-	expression tag	UNP P10104
A	1361	GLY	-	expression tag	UNP P10104
A	1362	SER	-	expression tag	UNP P10104
A	1363	ALA	-	expression tag	UNP P10104
A	1364	TRP	-	expression tag	UNP P10104
A	1365	SER	-	expression tag	UNP P10104
A	1366	HIS	-	expression tag	UNP P10104
A	1367	PRO	-	expression tag	UNP P10104
A	1368	GLN	-	expression tag	UNP P10104
A	1369	PHE	-	expression tag	UNP P10104
A	1370	GLU	-	expression tag	UNP P10104
A	1371	LYS	-	expression tag	UNP P10104
A	1372	SER	-	expression tag	UNP P10104
A	1373	ALA	-	expression tag	UNP P10104
B	885	PRO	ALA	engineered mutation	UNP A0A2R4KP93
B	962	PRO	SER	engineered mutation	UNP A0A2R4KP93
B	969	PRO	ALA	engineered mutation	UNP A0A2R4KP93
B	1012	PRO	ASN	engineered mutation	UNP A0A2R4KP93
B	1056	PRO	ALA	engineered mutation	UNP A0A2R4KP93
B	1057	PRO	VAL	engineered mutation	UNP A0A2R4KP93
B	1295	GLY	-	linker	UNP A0A2R4KP93

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1296	SER	-	linker	UNP A0A2R4KP93
B	1318	LEU	PHE	conflict	UNP P10104
B	1324	LEU	-	expression tag	UNP P10104
B	1325	GLU	-	expression tag	UNP P10104
B	1326	VAL	-	expression tag	UNP P10104
B	1327	LEU	-	expression tag	UNP P10104
B	1328	PHE	-	expression tag	UNP P10104
B	1329	GLN	-	expression tag	UNP P10104
B	1330	GLY	-	expression tag	UNP P10104
B	1331	PRO	-	expression tag	UNP P10104
B	1332	GLY	-	expression tag	UNP P10104
B	1333	HIS	-	expression tag	UNP P10104
B	1334	HIS	-	expression tag	UNP P10104
B	1335	HIS	-	expression tag	UNP P10104
B	1336	HIS	-	expression tag	UNP P10104
B	1337	HIS	-	expression tag	UNP P10104
B	1338	HIS	-	expression tag	UNP P10104
B	1339	HIS	-	expression tag	UNP P10104
B	1340	HIS	-	expression tag	UNP P10104
B	1341	SER	-	expression tag	UNP P10104
B	1342	ALA	-	expression tag	UNP P10104
B	1343	TRP	-	expression tag	UNP P10104
B	1344	SER	-	expression tag	UNP P10104
B	1345	HIS	-	expression tag	UNP P10104
B	1346	PRO	-	expression tag	UNP P10104
B	1347	GLN	-	expression tag	UNP P10104
B	1348	PHE	-	expression tag	UNP P10104
B	1349	GLU	-	expression tag	UNP P10104
B	1350	LYS	-	expression tag	UNP P10104
B	1351	GLY	-	expression tag	UNP P10104
B	1352	GLY	-	expression tag	UNP P10104
B	1353	GLY	-	expression tag	UNP P10104
B	1354	SER	-	expression tag	UNP P10104
B	1355	GLY	-	expression tag	UNP P10104
B	1356	GLY	-	expression tag	UNP P10104
B	1357	GLY	-	expression tag	UNP P10104
B	1358	GLY	-	expression tag	UNP P10104
B	1359	SER	-	expression tag	UNP P10104
B	1360	GLY	-	expression tag	UNP P10104
B	1361	GLY	-	expression tag	UNP P10104
B	1362	SER	-	expression tag	UNP P10104
B	1363	ALA	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1364	TRP	-	expression tag	UNP P10104
B	1365	SER	-	expression tag	UNP P10104
B	1366	HIS	-	expression tag	UNP P10104
B	1367	PRO	-	expression tag	UNP P10104
B	1368	GLN	-	expression tag	UNP P10104
B	1369	PHE	-	expression tag	UNP P10104
B	1370	GLU	-	expression tag	UNP P10104
B	1371	LYS	-	expression tag	UNP P10104
B	1372	SER	-	expression tag	UNP P10104
B	1373	ALA	-	expression tag	UNP P10104
C	885	PRO	ALA	engineered mutation	UNP A0A2R4KP93
C	962	PRO	SER	engineered mutation	UNP A0A2R4KP93
C	969	PRO	ALA	engineered mutation	UNP A0A2R4KP93
C	1012	PRO	ASN	engineered mutation	UNP A0A2R4KP93
C	1056	PRO	ALA	engineered mutation	UNP A0A2R4KP93
C	1057	PRO	VAL	engineered mutation	UNP A0A2R4KP93
C	1295	GLY	-	linker	UNP A0A2R4KP93
C	1296	SER	-	linker	UNP A0A2R4KP93
C	1318	LEU	PHE	conflict	UNP P10104
C	1324	LEU	-	expression tag	UNP P10104
C	1325	GLU	-	expression tag	UNP P10104
C	1326	VAL	-	expression tag	UNP P10104
C	1327	LEU	-	expression tag	UNP P10104
C	1328	PHE	-	expression tag	UNP P10104
C	1329	GLN	-	expression tag	UNP P10104
C	1330	GLY	-	expression tag	UNP P10104
C	1331	PRO	-	expression tag	UNP P10104
C	1332	GLY	-	expression tag	UNP P10104
C	1333	HIS	-	expression tag	UNP P10104
C	1334	HIS	-	expression tag	UNP P10104
C	1335	HIS	-	expression tag	UNP P10104
C	1336	HIS	-	expression tag	UNP P10104
C	1337	HIS	-	expression tag	UNP P10104
C	1338	HIS	-	expression tag	UNP P10104
C	1339	HIS	-	expression tag	UNP P10104
C	1340	HIS	-	expression tag	UNP P10104
C	1341	SER	-	expression tag	UNP P10104
C	1342	ALA	-	expression tag	UNP P10104
C	1343	TRP	-	expression tag	UNP P10104
C	1344	SER	-	expression tag	UNP P10104
C	1345	HIS	-	expression tag	UNP P10104
C	1346	PRO	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1347	GLN	-	expression tag	UNP P10104
C	1348	PHE	-	expression tag	UNP P10104
C	1349	GLU	-	expression tag	UNP P10104
C	1350	LYS	-	expression tag	UNP P10104
C	1351	GLY	-	expression tag	UNP P10104
C	1352	GLY	-	expression tag	UNP P10104
C	1353	GLY	-	expression tag	UNP P10104
C	1354	SER	-	expression tag	UNP P10104
C	1355	GLY	-	expression tag	UNP P10104
C	1356	GLY	-	expression tag	UNP P10104
C	1357	GLY	-	expression tag	UNP P10104
C	1358	GLY	-	expression tag	UNP P10104
C	1359	SER	-	expression tag	UNP P10104
C	1360	GLY	-	expression tag	UNP P10104
C	1361	GLY	-	expression tag	UNP P10104
C	1362	SER	-	expression tag	UNP P10104
C	1363	ALA	-	expression tag	UNP P10104
C	1364	TRP	-	expression tag	UNP P10104
C	1365	SER	-	expression tag	UNP P10104
C	1366	HIS	-	expression tag	UNP P10104
C	1367	PRO	-	expression tag	UNP P10104
C	1368	GLN	-	expression tag	UNP P10104
C	1369	PHE	-	expression tag	UNP P10104
C	1370	GLU	-	expression tag	UNP P10104
C	1371	LYS	-	expression tag	UNP P10104
C	1372	SER	-	expression tag	UNP P10104
C	1373	ALA	-	expression tag	UNP P10104

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



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

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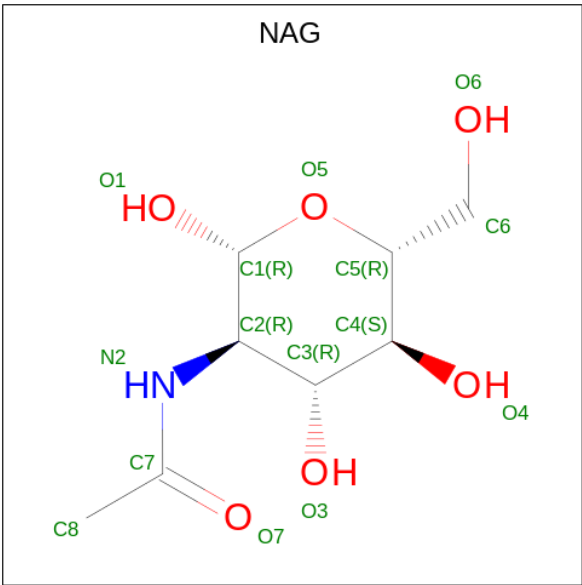
Mol	Chain	Residues	Atoms				AltConf	Trace
2	I	2	Total	C	N	O	0	0
			28	16	2	10		
2	J	2	Total	C	N	O	0	0
			28	16	2	10		
2	K	2	Total	C	N	O	0	0
			28	16	2	10		
2	L	2	Total	C	N	O	0	0
			28	16	2	10		
2	N	2	Total	C	N	O	0	0
			28	16	2	10		
2	O	2	Total	C	N	O	0	0
			28	16	2	10		

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



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

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



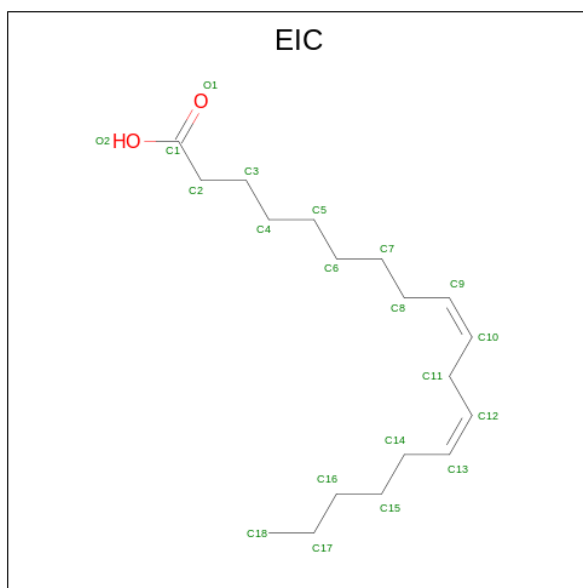
Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4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

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

- Molecule 5 is LINOLEIC ACID (CCD ID: EIC) (formula: $C_{18}H_{32}O_2$) (labeled as "Ligand of Interest" by depositor).

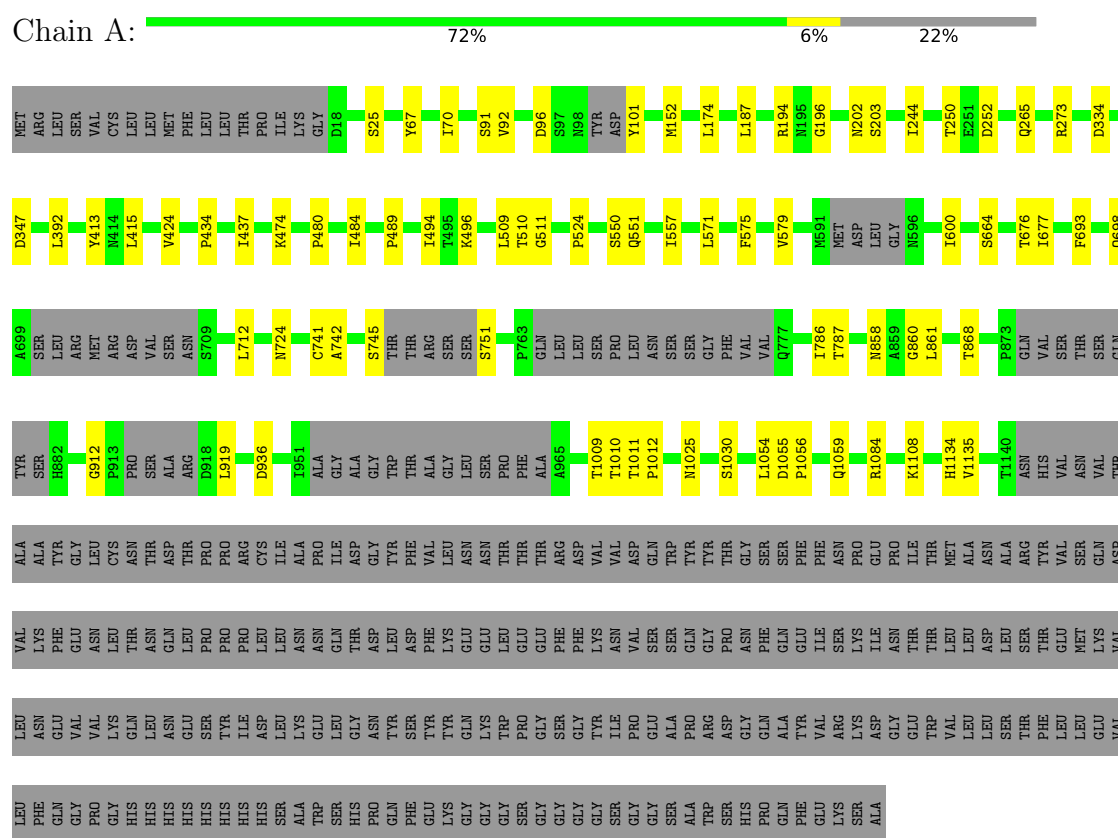


Mol	Chain	Residues	Atoms			AltConf
5	A	1	Total	C	O	0
			20	18	2	
5	B	1	Total	C	O	0
			20	18	2	
5	C	1	Total	C	O	0
			20	18	2	

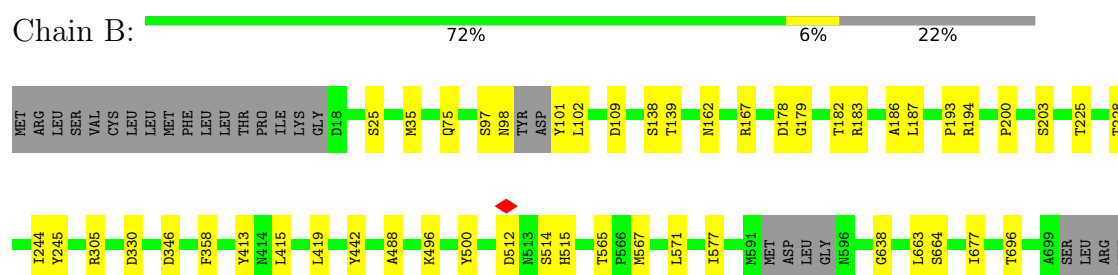
3 Residue-property plots

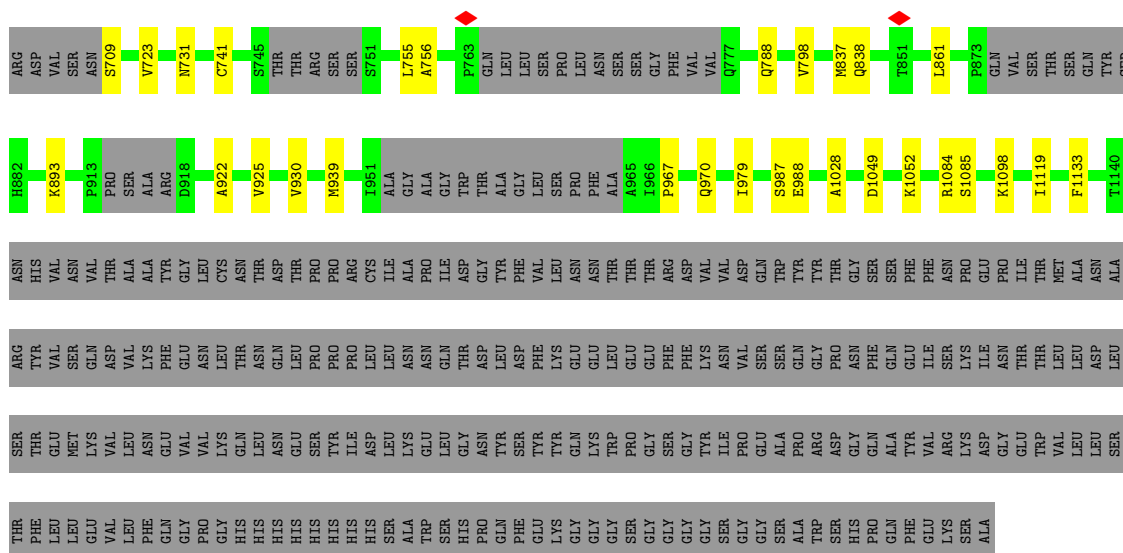
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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,Fibritin

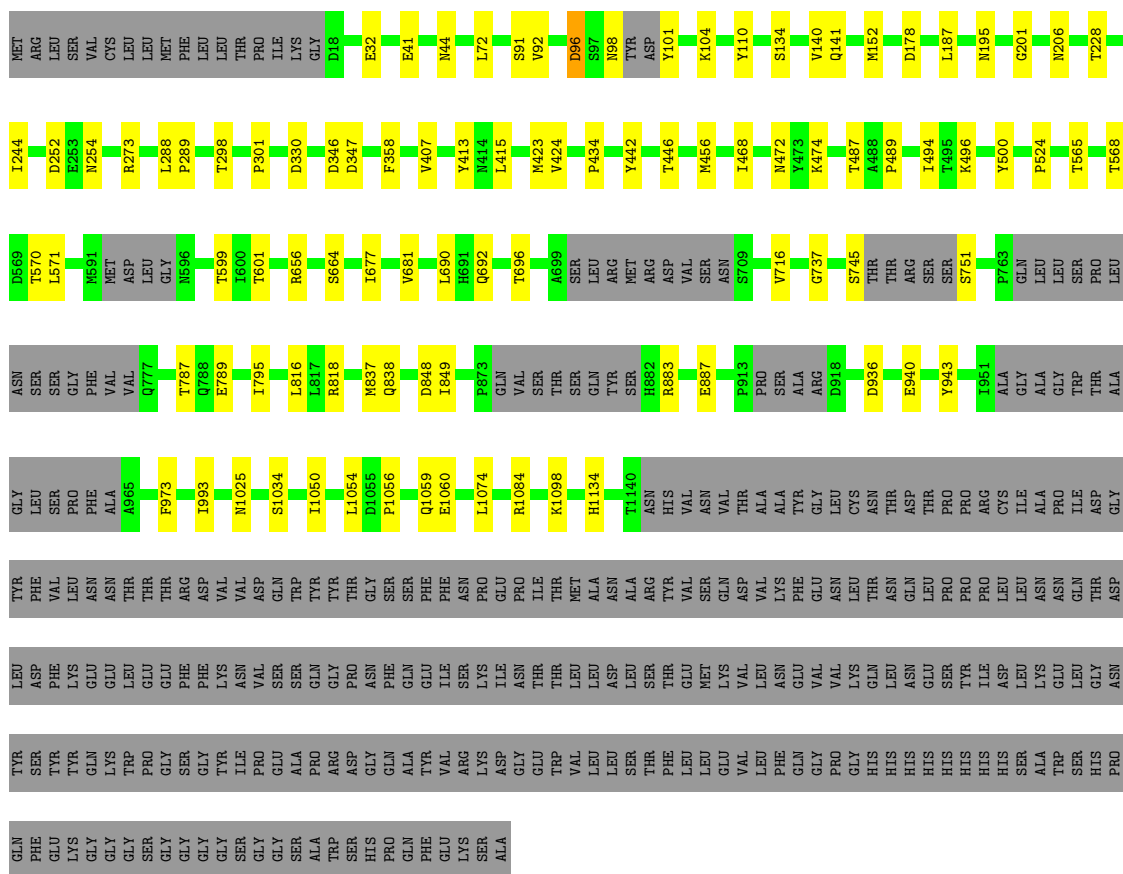


• Molecule 1: Spike glycoprotein,Fibritin





- Molecule 1: Spike glycoprotein, Fibrin



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

Chain D:  100%



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

Chain F:  100%



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

Chain G:  100%



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

Chain I:  100%



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

Chain J:  50% 50%



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

Chain K:  100%



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

Chain L:  100%



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

Chain N:  50% 50%



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

Chain O:  100%



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

Chain E:  100%



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

Chain H:  100%



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

Chain M:  67% 33%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	111480	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.248	Depositor
Minimum map value	-0.122	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.022	Depositor
Map size (\AA)	422.40002, 422.40002, 422.40002	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.32, 1.32, 1.32	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: NAG, EIC, BMA

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	0.17	0/8414	0.36	0/11427
1	B	0.17	0/8414	0.35	0/11427
1	C	0.19	0/8414	0.37	0/11427
All	All	0.18	0/25242	0.36	0/34281

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	8228	0	7916	52	0
1	B	8228	0	7916	48	0
1	C	8228	0	7916	56	0
2	D	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	I	28	0	25	0	0
2	J	28	0	25	0	0
2	K	28	0	25	0	0
2	L	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	N	28	0	25	0	0
2	O	28	0	25	0	0
3	E	39	0	34	0	0
3	H	39	0	34	0	0
3	M	39	0	34	1	0
4	A	98	0	91	1	0
4	B	98	0	91	2	0
4	C	98	0	91	0	0
5	A	20	0	31	2	0
5	B	20	0	31	2	0
5	C	20	0	31	3	0
All	All	25407	0	24441	152	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 3.

All (152) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:745:SER:HG	1:A:751:SER:N	1.84	0.76
1:A:1009:THR:O	1:A:1010:THR:HG22	1.94	0.68
1:B:98:ASN:O	1:B:101:TYR:N	2.29	0.65
1:B:861:LEU:HD13	4:B:1404:NAG:H82	1.77	0.65
1:A:496:LYS:HG2	1:A:571:LEU:HB2	1.81	0.63
1:B:638:GLY:HA2	1:C:72:LEU:HD21	1.81	0.62
1:C:664:SER:HB2	1:C:677:ILE:HB	1.82	0.62
1:B:162:ASN:HD21	1:B:203:SER:HA	1.66	0.60
1:C:252:ASP:HB2	1:C:273:ARG:HH22	1.66	0.60
5:B:1408:EIC:H31	1:C:424:VAL:HG11	1.85	0.59
1:C:32:GLU:HG2	1:C:104:LYS:HD2	1.85	0.59
1:B:98:ASN:C	1:B:101:TYR:N	2.61	0.58
1:B:664:SER:HB2	1:B:677:ILE:HB	1.84	0.58
1:A:489:PRO:HD2	1:A:494:ILE:HD12	1.86	0.58
1:B:496:LYS:HG2	1:B:571:LEU:HB2	1.86	0.58
1:B:97:SER:O	1:B:101:TYR:N	2.37	0.58
1:C:496:LYS:HG2	1:C:571:LEU:HB2	1.86	0.57
1:A:912:GLY:O	1:C:656:ARG:NH1	2.37	0.57
1:A:413:TYR:HE2	1:A:415:LEU:HD13	1.71	0.55
1:C:254:ASN:HA	3:M:1:NAG:HN2	1.71	0.55
1:A:265:GLN:HA	1:C:407:VAL:HG21	1.88	0.55
1:A:202:ASN:OD1	1:A:203:SER:N	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:193:PRO:HB3	1:B:200:PRO:HG2	1.88	0.55
1:B:514:SER:OG	1:B:515:HIS:N	2.40	0.55
1:A:509:LEU:HB2	1:A:557:ILE:HB	1.89	0.54
1:B:500:TYR:HB2	1:B:565:THR:HB	1.90	0.54
1:A:187:LEU:HD23	1:A:244:ILE:HG12	1.89	0.54
1:A:1011:THR:HB	1:A:1012:PRO:HD3	1.89	0.54
1:C:500:TYR:HB2	1:C:565:THR:HB	1.90	0.54
1:C:178:ASP:OD2	1:C:228:THR:N	2.39	0.54
1:B:514:SER:O	1:B:515:HIS:ND1	2.41	0.53
1:C:187:LEU:HD23	1:C:244:ILE:HG12	1.89	0.53
1:A:860:GLY:HA2	1:A:868:THR:HB	1.91	0.53
1:C:1025:ASN:OD1	1:C:1084:ARG:NH2	2.41	0.53
1:A:936:ASP:OD1	1:A:936:ASP:N	2.41	0.53
1:C:1056:PRO:HA	1:C:1059:GLN:HB2	1.90	0.53
1:C:134:SER:HA	1:C:141:GLN:HA	1.90	0.53
1:A:91:SER:OG	1:A:92:VAL:N	2.42	0.53
1:B:138:SER:OG	1:B:139:THR:N	2.42	0.52
1:A:437:ILE:HG21	5:C:1401:EIC:H91	1.92	0.52
1:B:186:ALA:HB3	1:B:245:TYR:HB2	1.92	0.52
1:B:837:MET:HE1	1:B:1085:SER:HB3	1.91	0.52
1:C:489:PRO:HD2	1:C:494:ILE:HD12	1.92	0.51
1:C:936:ASP:N	1:C:936:ASP:OD1	2.41	0.51
1:C:795:ILE:O	1:C:838:GLN:NE2	2.37	0.51
1:A:424:VAL:HG11	5:C:1401:EIC:H21	1.92	0.51
1:B:346:ASP:HB3	1:B:696:THR:HB	1.92	0.51
1:A:1055:ASP:OD1	1:A:1055:ASP:N	2.35	0.50
1:B:75:GLN:OE1	1:C:818:ARG:NE	2.45	0.50
1:A:787:THR:HG22	1:A:1134:HIS:HB2	1.94	0.50
1:C:789:GLU:OE2	1:C:1098:LYS:NZ	2.41	0.50
1:C:206:ASN:OD1	1:C:206:ASN:N	2.45	0.50
1:A:152:MET:HG3	1:A:174:LEU:HD23	1.94	0.50
1:A:741:CYS:SG	1:A:742:ALA:N	2.85	0.50
1:C:716:VAL:HG11	1:C:737:GLY:HA3	1.94	0.50
1:B:25:SER:O	1:B:194:ARG:NH2	2.45	0.50
5:B:1408:EIC:H52	1:C:434:PRO:HB3	1.93	0.49
1:A:1056:PRO:HA	1:A:1059:GLN:HB2	1.95	0.49
1:A:1025:ASN:OD1	1:A:1084:ARG:NH2	2.43	0.49
1:A:434:PRO:HB3	5:C:1401:EIC:H52	1.95	0.48
1:C:745:SER:HG	1:C:751:SER:N	2.10	0.48
1:B:179:GLY:O	1:B:182:THR:OG1	2.30	0.48
1:B:1028:ALA:HB2	1:B:1084:ARG:HH12	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:480:PRO:HG2	1:A:579:VAL:HG23	1.95	0.48
1:B:512:ASP:OD1	1:B:512:ASP:N	2.47	0.48
1:C:787:THR:HG23	1:C:1134:HIS:HB2	1.96	0.48
1:A:510:THR:OG1	1:A:511:GLY:N	2.47	0.47
1:C:347:ASP:OD1	1:C:347:ASP:N	2.39	0.47
1:C:1054:LEU:O	1:C:1059:GLN:NE2	2.47	0.47
1:A:196:GLY:H	1:A:202:ASN:HB2	1.79	0.47
1:B:1049:ASP:HA	1:B:1052:LYS:HG2	1.97	0.47
1:A:25:SER:O	1:A:194:ARG:NH1	2.45	0.47
1:B:755:LEU:HD21	4:B:1403:NAG:H82	1.97	0.46
5:A:1408:EIC:H62	1:B:419:LEU:HD21	1.96	0.46
1:B:183:ARG:HD2	1:B:225:THR:HG23	1.98	0.46
1:C:407:VAL:HG22	1:C:446:THR:HG23	1.98	0.46
1:C:468:ILE:HA	1:C:472:ASN:HD22	1.80	0.45
1:C:690:LEU:HG	1:C:692:GLN:HG2	1.98	0.45
1:B:731:ASN:OD1	1:B:731:ASN:N	2.49	0.45
1:A:861:LEU:HD21	4:A:1405:NAG:H82	1.98	0.45
1:B:567:MET:HE2	1:B:567:MET:HB2	1.86	0.45
1:A:664:SER:HB2	1:A:677:ILE:HB	1.97	0.45
1:A:919:LEU:HD11	1:A:1030:SER:HB2	1.98	0.45
1:C:837:MET:HE3	1:C:837:MET:HB3	1.88	0.45
1:C:940:GLU:HA	1:C:943:TYR:HD1	1.82	0.45
1:A:434:PRO:HA	1:A:437:ILE:HD12	1.98	0.45
1:B:741:CYS:N	1:B:756:ALA:O	2.50	0.45
1:A:1108:LYS:HD3	1:A:1108:LYS:HA	1.78	0.44
1:C:41:GLU:HA	1:C:44:ASN:HD21	1.82	0.44
1:A:550:SER:OG	1:A:551:GLN:N	2.50	0.44
1:B:798:VAL:HG22	1:B:930:VAL:HG22	1.99	0.44
1:A:1054:LEU:O	1:A:1059:GLN:NE2	2.50	0.44
1:C:1098:LYS:HE3	1:C:1098:LYS:HB2	1.83	0.44
1:C:423:MET:HB3	1:C:487:THR:HB	2.00	0.44
1:B:187:LEU:HD23	1:B:244:ILE:HD12	2.00	0.43
1:C:848:ASP:OD1	1:C:848:ASP:N	2.51	0.43
1:A:250:THR:O	1:A:273:ARG:NH2	2.51	0.43
1:B:413:TYR:HE2	1:B:415:LEU:HD13	1.83	0.43
1:A:474:LYS:HD3	1:A:524:PRO:HG3	2.00	0.43
1:B:939:MET:HE2	1:B:939:MET:HB3	1.89	0.43
1:C:973:PHE:HE1	1:C:993:ILE:HG13	1.81	0.43
1:B:35:MET:HE1	1:B:109:ASP:HA	1.99	0.43
1:C:346:ASP:HB3	1:C:696:THR:HB	2.00	0.43
1:B:922:ALA:HA	1:B:925:VAL:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:788:GLN:HE21	1:B:1133:PHE:HE1	1.67	0.43
1:A:474:LYS:HB3	1:A:524:PRO:HA	2.01	0.42
1:C:152:MET:HB3	1:C:298:THR:HG22	2.00	0.42
1:A:712:LEU:HD21	1:A:724:ASN:HB2	2.00	0.42
1:A:858:ASN:OD1	1:A:858:ASN:N	2.48	0.42
1:C:474:LYS:HB3	1:C:524:PRO:HA	2.01	0.42
1:C:1074:LEU:HD23	1:C:1074:LEU:HA	1.90	0.42
1:C:883:ARG:HG2	1:C:887:GLU:HB3	2.00	0.42
1:B:178:ASP:OD2	1:B:228:THR:N	2.51	0.42
1:A:1010:THR:O	1:A:1010:THR:HG23	2.19	0.42
1:B:979:ILE:HD13	1:B:1119:ILE:HD11	2.01	0.42
1:C:195:ASN:HA	1:C:201:GLY:HA3	2.01	0.42
1:A:67:TYR:HB3	1:A:70:ILE:HD11	2.02	0.41
1:A:252:ASP:HB2	1:A:273:ARG:HH12	1.86	0.41
1:B:488:ALA:HB3	1:B:571:LEU:HB3	2.01	0.41
1:B:967:PRO:HD2	1:B:970:GLN:HG3	2.02	0.41
1:B:987:SER:OG	1:B:988:GLU:OE2	2.31	0.41
1:A:484:ILE:HB	1:A:575:PHE:HB2	2.01	0.41
1:A:600:ILE:HD12	1:A:600:ILE:HA	1.96	0.41
1:A:693:PHE:O	1:A:698:GLN:NE2	2.52	0.41
1:C:330:ASP:HB2	1:C:358:PHE:CE1	2.55	0.41
1:C:568:THR:OG1	1:C:570:THR:O	2.38	0.41
1:A:1054:LEU:HD23	1:A:1054:LEU:HA	1.94	0.41
5:A:1408:EIC:H82	1:B:577:ILE:HD13	2.02	0.41
1:B:893:LYS:HE3	1:B:893:LYS:HB3	1.90	0.41
1:C:96:ASP:OD2	1:C:101:TYR:HD2	2.03	0.41
1:C:599:THR:HG22	1:C:601:THR:H	1.86	0.41
1:C:690:LEU:HD12	1:C:690:LEU:HA	1.87	0.41
1:C:1050:ILE:HG23	1:C:1054:LEU:HD12	2.02	0.41
1:B:663:LEU:HD23	1:B:663:LEU:HA	1.95	0.41
1:C:288:LEU:HA	1:C:289:PRO:HD3	1.92	0.41
1:C:456:MET:HE3	1:C:456:MET:HB3	1.98	0.41
1:A:96:ASP:OD2	1:A:101:TYR:N	2.53	0.41
1:A:919:LEU:HD23	1:A:919:LEU:HA	1.94	0.41
1:B:167:ARG:HA	1:B:167:ARG:HD2	1.75	0.41
1:A:392:LEU:HD11	1:A:575:PHE:HE1	1.86	0.41
1:C:413:TYR:OH	1:C:442:TYR:OH	2.29	0.41
1:A:334:ASP:OD1	1:A:334:ASP:N	2.46	0.40
1:B:709:SER:HB2	1:B:723:VAL:HG12	2.03	0.40
1:C:413:TYR:HE2	1:C:415:LEU:HD13	1.85	0.40
1:A:347:ASP:OD2	1:A:676:THR:OG1	2.31	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:91:SER:OG	1:C:92:VAL:N	2.54	0.40
1:B:102:LEU:HD13	1:B:305:ARG:HB3	2.03	0.40
1:B:330:ASP:HB2	1:B:358:PHE:CE1	2.56	0.40
1:B:1098:LYS:HE2	1:B:1098:LYS:HB2	1.99	0.40
1:C:110:TYR:HB2	1:C:301:PRO:HG3	2.04	0.40
1:C:816:LEU:HD11	1:C:1060:GLU:HG3	2.03	0.40
1:A:786:ILE:HG12	1:A:1135:VAL:HG22	2.02	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	1047/1373 (76%)	1017 (97%)	30 (3%)	0	100	100
1	B	1047/1373 (76%)	1011 (97%)	36 (3%)	0	100	100
1	C	1047/1373 (76%)	1009 (96%)	38 (4%)	0	100	100
All	All	3141/4119 (76%)	3037 (97%)	104 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	919/1186 (78%)	919 (100%)	0	100	100
1	B	919/1186 (78%)	917 (100%)	2 (0%)	92	96
1	C	919/1186 (78%)	913 (99%)	6 (1%)	81	90
All	All	2757/3558 (78%)	2749 (100%)	8 (0%)	90	95

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	442	TYR
1	B	838	GLN
1	C	96	ASP
1	C	98	ASN
1	C	140	VAL
1	C	681	VAL
1	C	849	ILE
1	C	1034	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	HIS
1	A	118	ASN
1	A	141	GLN
1	A	373	HIS
1	A	685	HIS
1	A	692	GLN
1	A	698	GLN
1	A	1068	ASN
1	A	1080	GLN
1	B	118	ASN
1	B	162	ASN
1	B	171	HIS
1	B	265	GLN
1	B	350	GLN
1	B	425	ASN
1	B	532	ASN
1	B	631	GLN
1	B	641	ASN
1	B	733	GLN
1	B	796	GLN
1	B	804	GLN

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Mol	Chain	Res	Type
1	B	983	GLN
1	B	1068	ASN
1	B	1106	GLN
1	B	1125	ASN
1	C	81	HIS
1	C	239	GLN
1	C	240	ASN
1	C	345	HIS
1	C	381	GLN
1	C	425	ASN
1	C	491	ASN
1	C	762	GLN
1	C	882	HIS
1	C	1027	GLN
1	C	1081	GLN

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](#)

27 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	0.25	0	17,19,21	0.52	0
2	NAG	D	2	2	14,14,15	0.26	0	17,19,21	0.44	0
3	NAG	E	1	1,3	14,14,15	0.28	0	17,19,21	0.57	0
3	NAG	E	2	3	14,14,15	0.25	0	17,19,21	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BMA	E	3	3	11,11,12	0.58	0	15,15,17	0.89	0
2	NAG	F	1	2,1	14,14,15	0.21	0	17,19,21	0.38	0
2	NAG	F	2	2	14,14,15	0.24	0	17,19,21	0.52	0
2	NAG	G	1	2,1	14,14,15	0.31	0	17,19,21	0.56	0
2	NAG	G	2	2	14,14,15	0.27	0	17,19,21	0.42	0
3	NAG	H	1	1,3	14,14,15	0.27	0	17,19,21	0.57	0
3	NAG	H	2	3	14,14,15	0.22	0	17,19,21	0.44	0
3	BMA	H	3	3	11,11,12	0.61	0	15,15,17	0.88	0
2	NAG	I	1	2,1	14,14,15	0.26	0	17,19,21	0.44	0
2	NAG	I	2	2	14,14,15	0.19	0	17,19,21	0.47	0
2	NAG	J	1	2,1	14,14,15	0.30	0	17,19,21	0.67	1 (5%)
2	NAG	J	2	2	14,14,15	0.29	0	17,19,21	0.42	0
2	NAG	K	1	2,1	14,14,15	0.22	0	17,19,21	0.60	0
2	NAG	K	2	2	14,14,15	0.25	0	17,19,21	0.46	0
2	NAG	L	1	2,1	14,14,15	0.24	0	17,19,21	0.56	0
2	NAG	L	2	2	14,14,15	0.26	0	17,19,21	0.48	0
3	NAG	M	1	1,3	14,14,15	0.27	0	17,19,21	0.53	0
3	NAG	M	2	3	14,14,15	0.23	0	17,19,21	0.41	0
3	BMA	M	3	3	11,11,12	0.58	0	15,15,17	0.89	0
2	NAG	N	1	2,1	14,14,15	0.34	0	17,19,21	0.91	2 (11%)
2	NAG	N	2	2	14,14,15	0.27	0	17,19,21	0.52	0
2	NAG	O	1	2,1	14,14,15	0.22	0	17,19,21	0.41	0
2	NAG	O	2	2	14,14,15	0.24	0	17,19,21	0.45	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
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
3	NAG	E	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
2	NAG	F	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	3/6/23/26	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
3	NAG	H	1	1,3	-	1/6/23/26	0/1/1/1
3	NAG	H	2	3	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BMA	H	3	3	-	0/2/19/22	0/1/1/1
2	NAG	I	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	I	2	2	-	2/6/23/26	0/1/1/1
2	NAG	J	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	J	2	2	-	1/6/23/26	0/1/1/1
2	NAG	K	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	K	2	2	-	0/6/23/26	0/1/1/1
2	NAG	L	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	L	2	2	-	0/6/23/26	0/1/1/1
3	NAG	M	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	M	2	3	-	2/6/23/26	0/1/1/1
3	BMA	M	3	3	-	0/2/19/22	0/1/1/1
2	NAG	N	1	2,1	-	1/6/23/26	0/1/1/1
2	NAG	N	2	2	-	0/6/23/26	0/1/1/1
2	NAG	O	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	O	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	1	NAG	C2-N2-C7	2.43	126.36	122.90
2	J	1	NAG	C1-O5-C5	2.30	115.31	112.19
2	N	1	NAG	C1-O5-C5	2.02	114.93	112.19

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	2	NAG	O5-C5-C6-O6
3	M	2	NAG	O5-C5-C6-O6
3	M	2	NAG	C4-C5-C6-O6
2	O	1	NAG	O5-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	O	1	NAG	C4-C5-C6-O6
2	I	1	NAG	O5-C5-C6-O6
2	F	2	NAG	C8-C7-N2-C2
2	F	2	NAG	O7-C7-N2-C2
2	I	2	NAG	C8-C7-N2-C2
2	I	2	NAG	O7-C7-N2-C2

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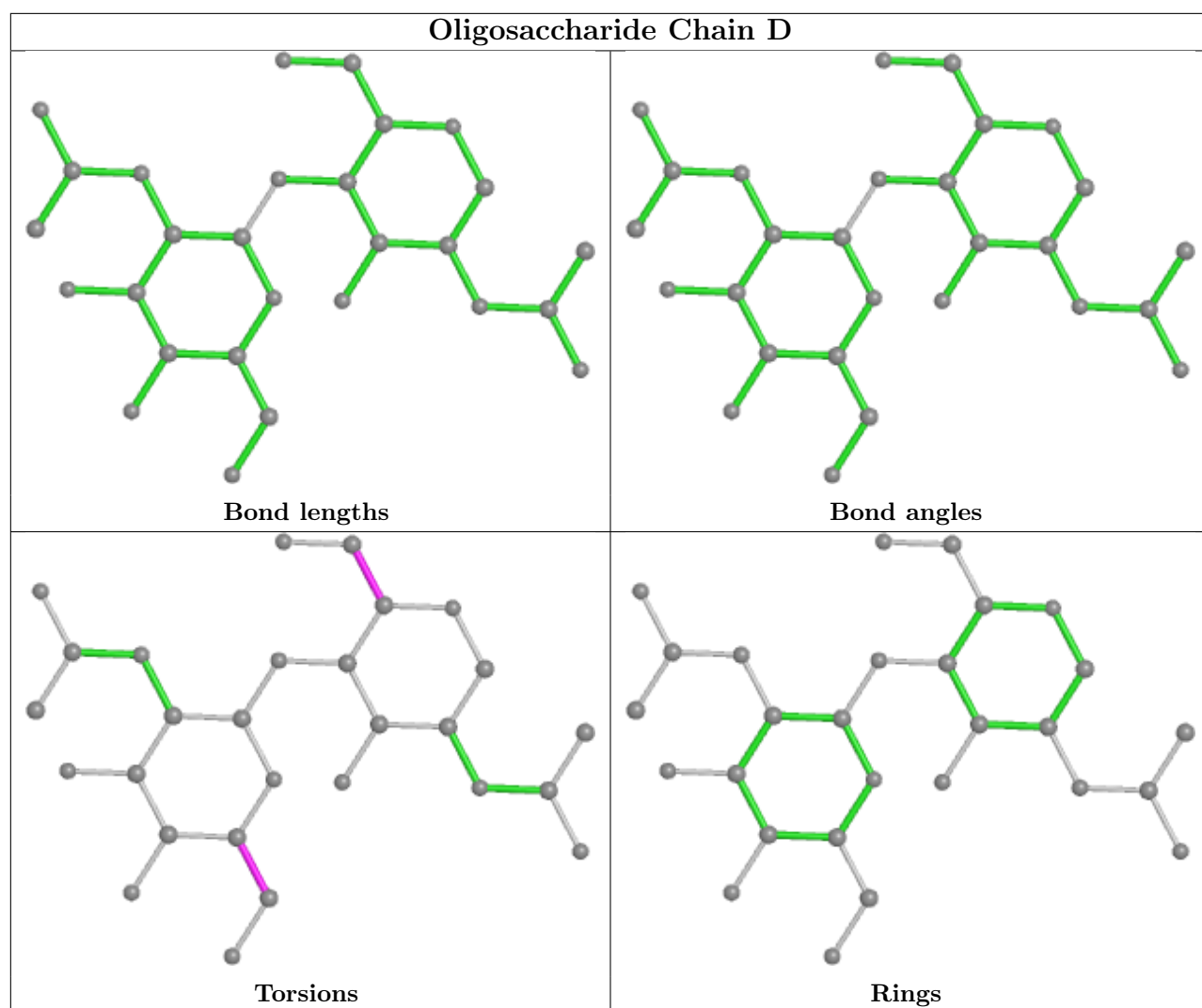
Mol	Chain	Res	Type	Atoms
2	K	1	NAG	C8-C7-N2-C2
2	K	1	NAG	O7-C7-N2-C2
3	E	1	NAG	C8-C7-N2-C2
3	E	1	NAG	O7-C7-N2-C2
2	L	1	NAG	O5-C5-C6-O6
3	E	2	NAG	O5-C5-C6-O6
2	I	1	NAG	C4-C5-C6-O6
2	K	1	NAG	O5-C5-C6-O6
2	D	1	NAG	O5-C5-C6-O6
2	L	1	NAG	C4-C5-C6-O6
2	K	1	NAG	C4-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	N	1	NAG	C3-C2-N2-C7
2	F	1	NAG	O5-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6

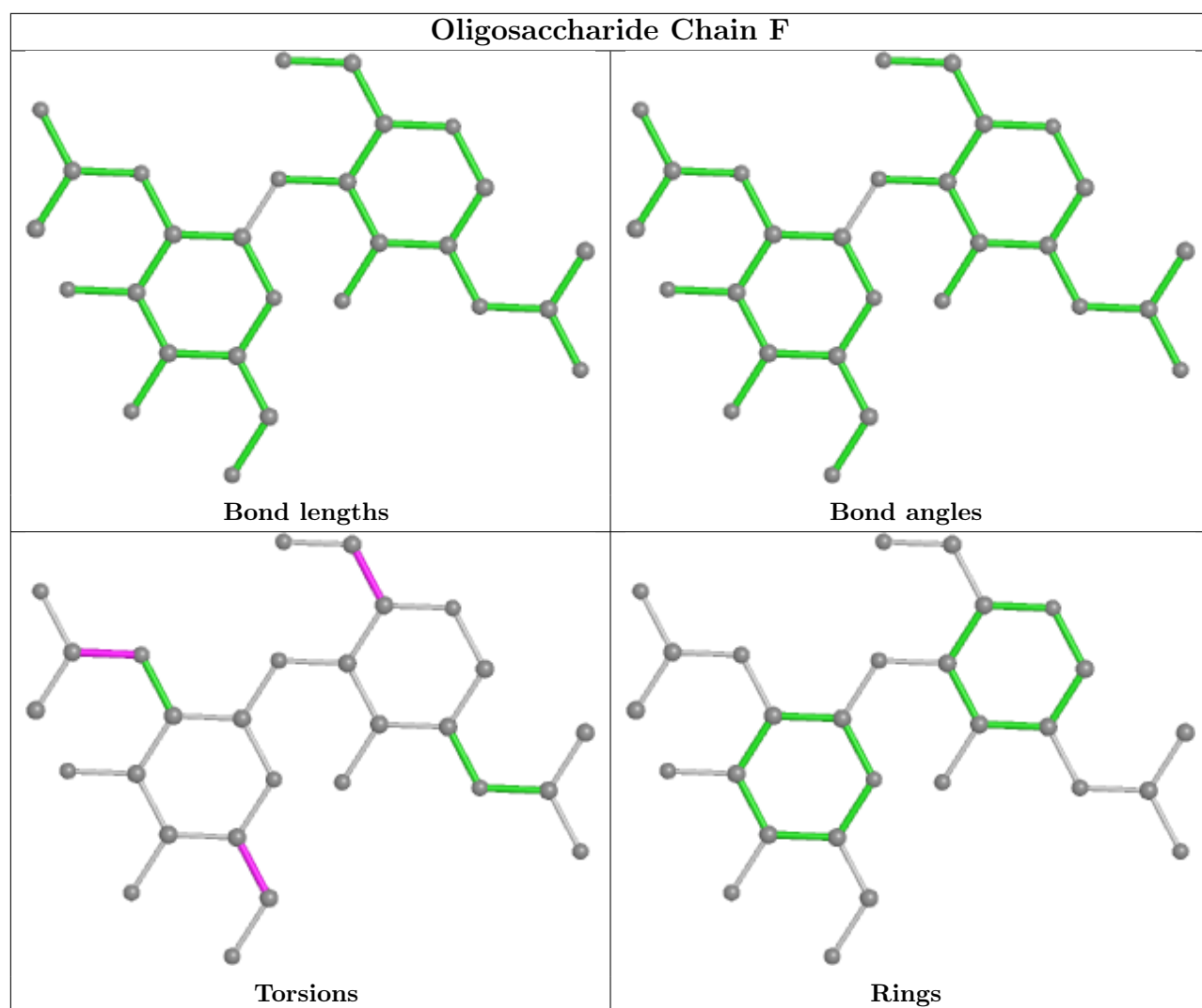
There are no ring outliers.

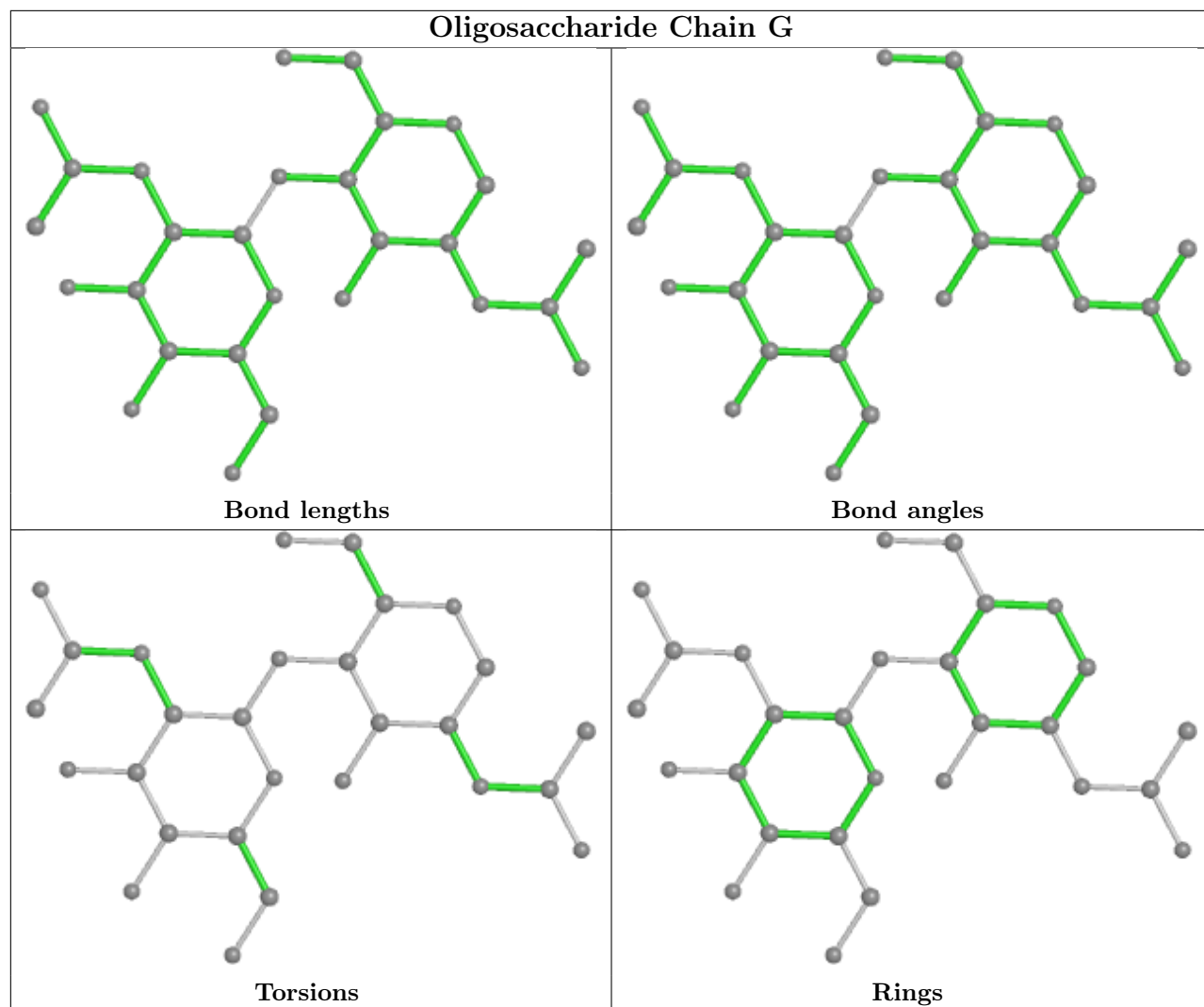
1 monomer is involved in 1 short contact:

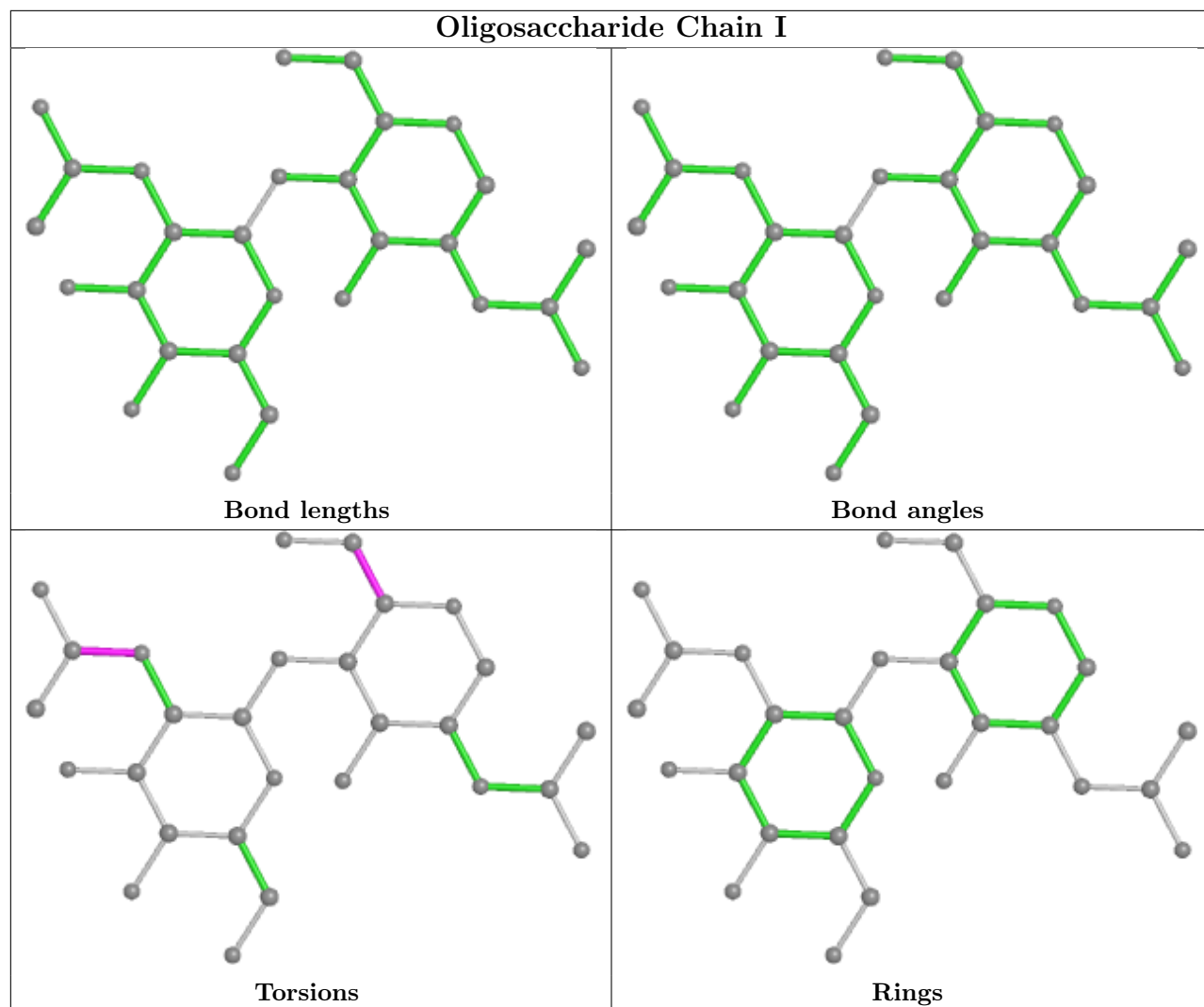
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	M	1	NAG	1	0

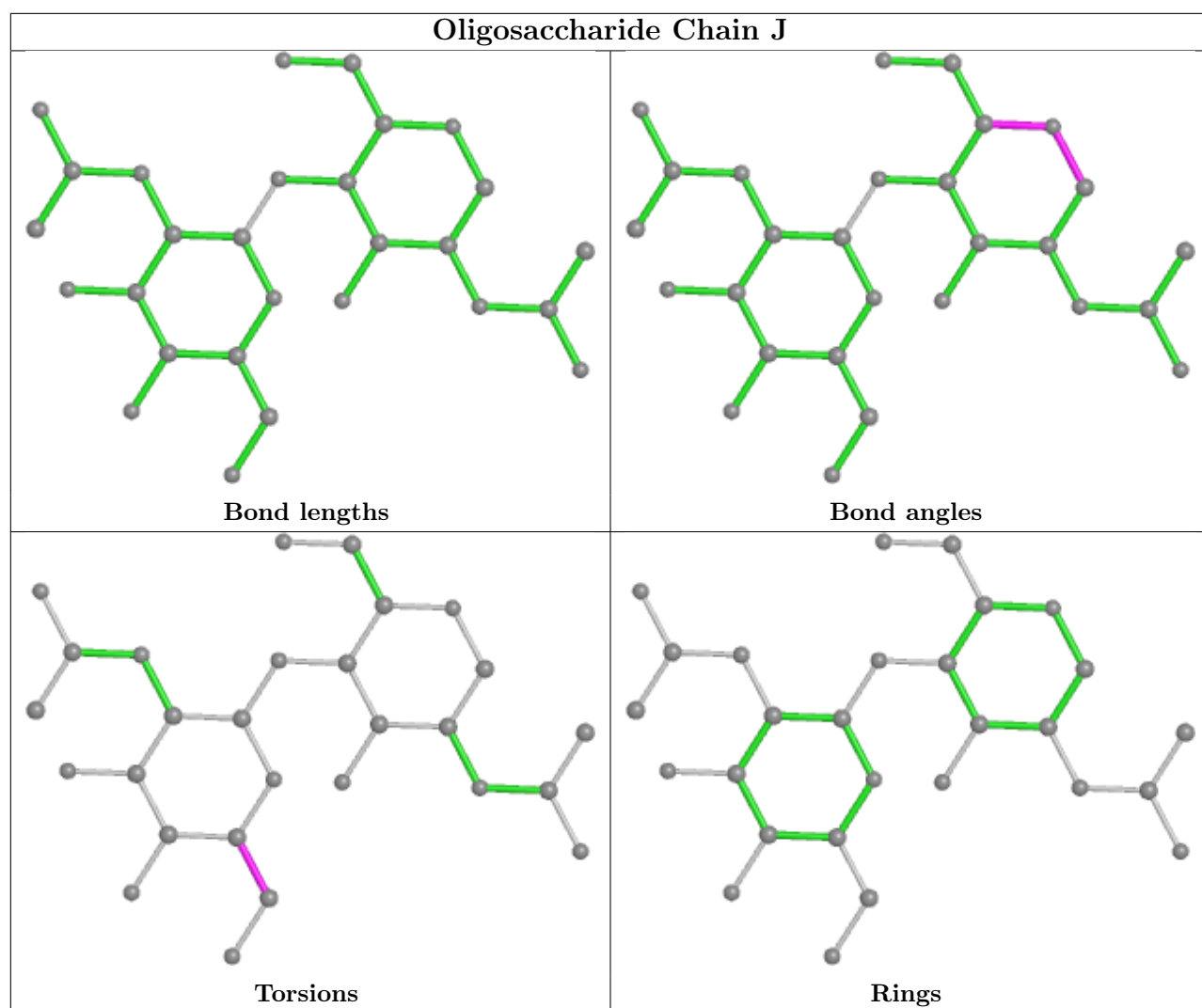
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

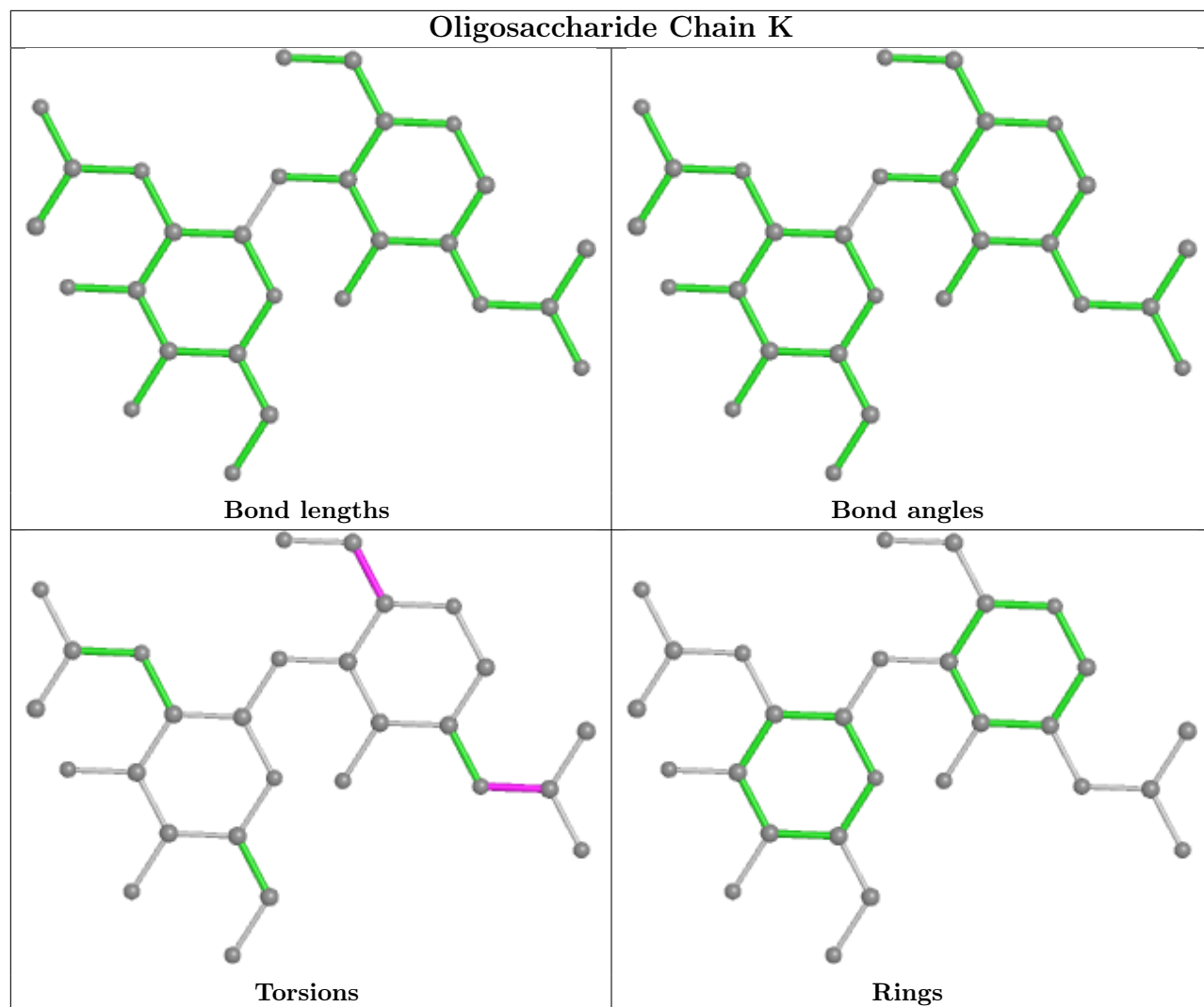


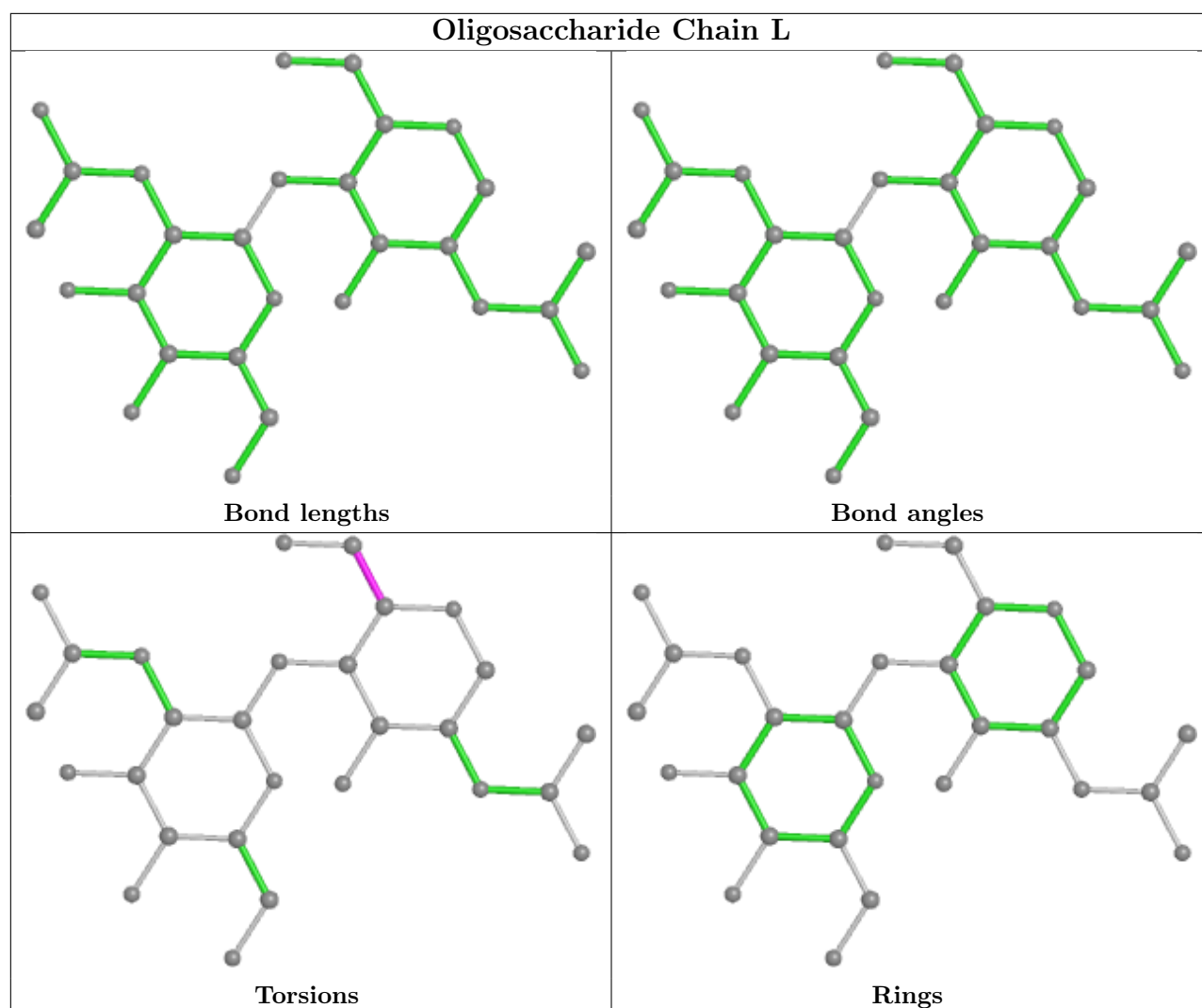


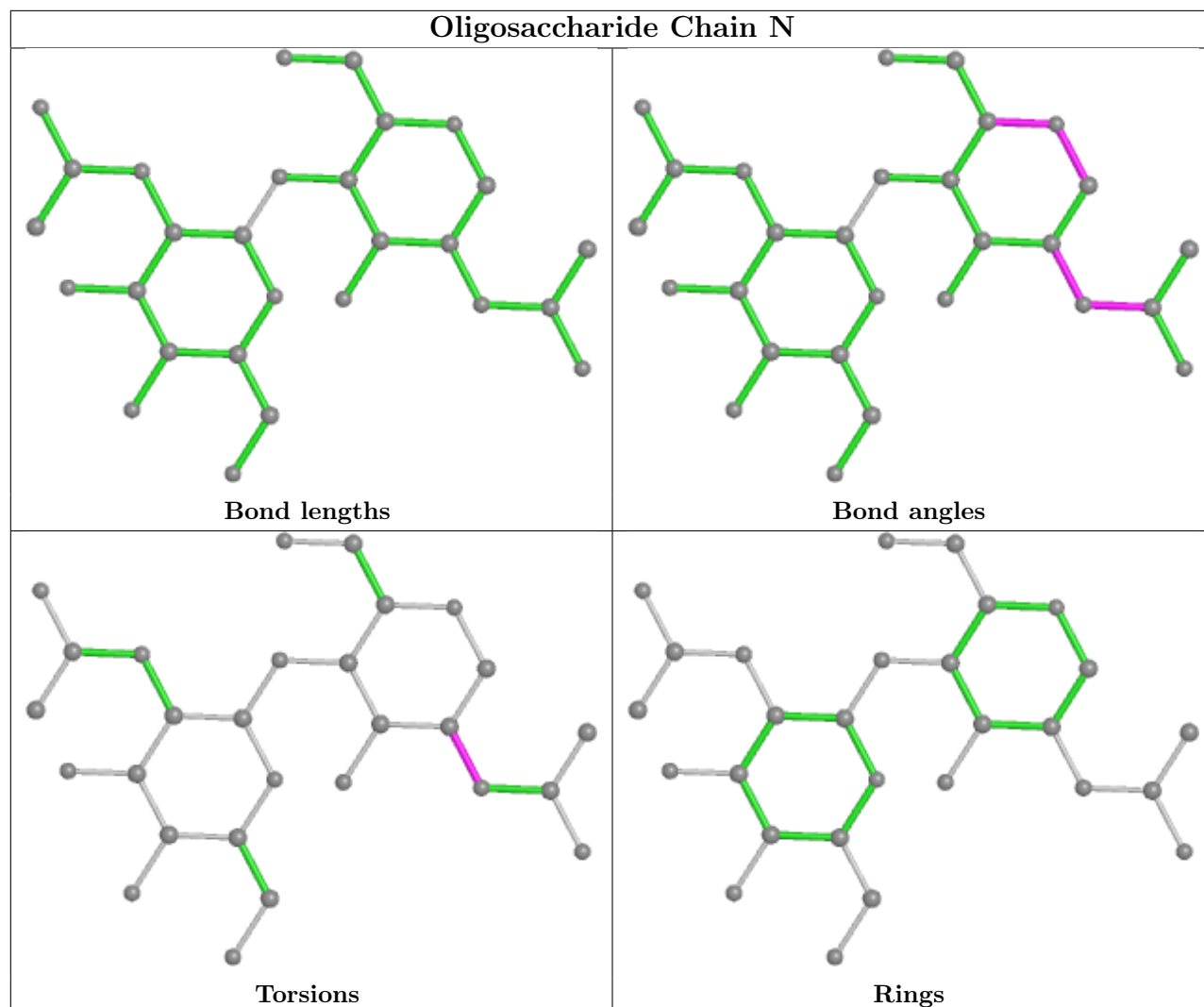


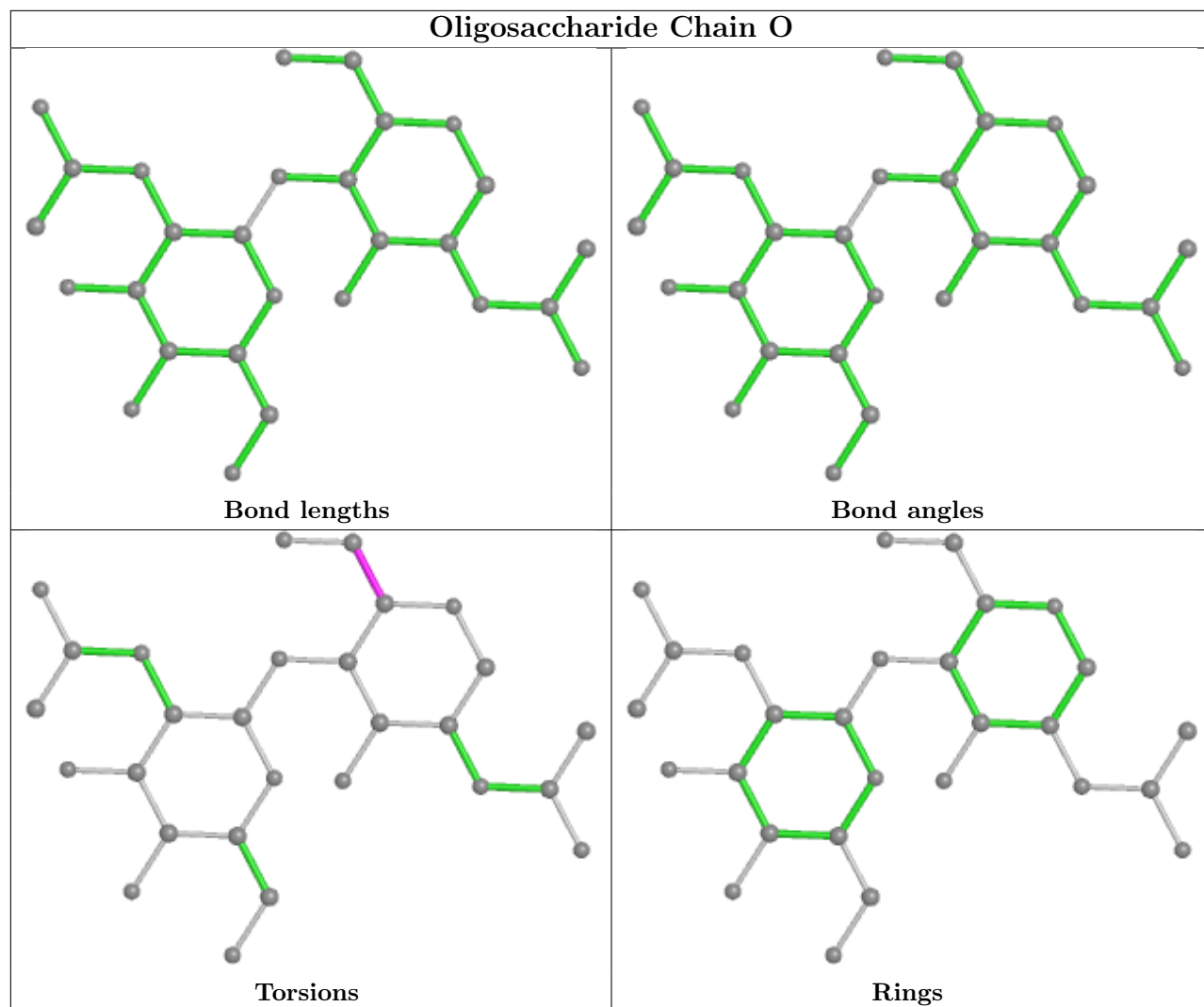


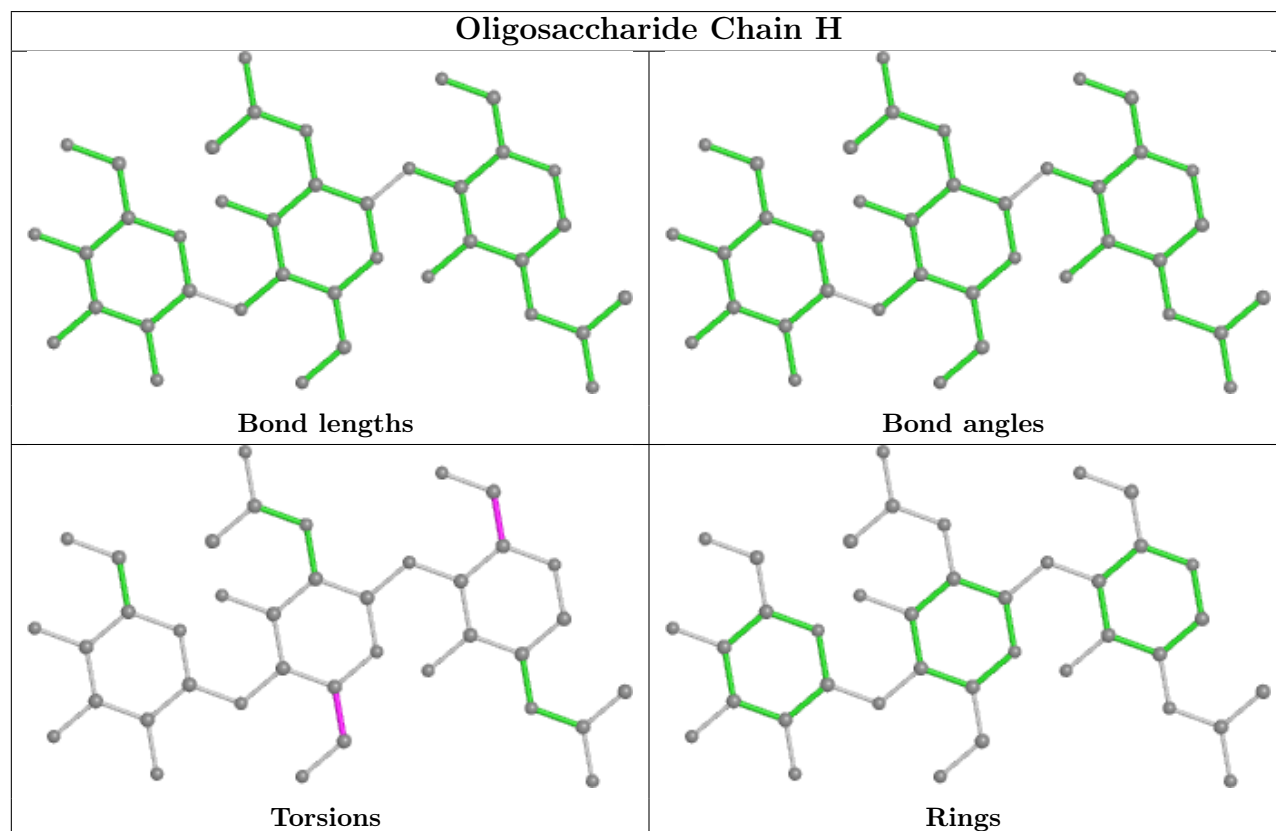
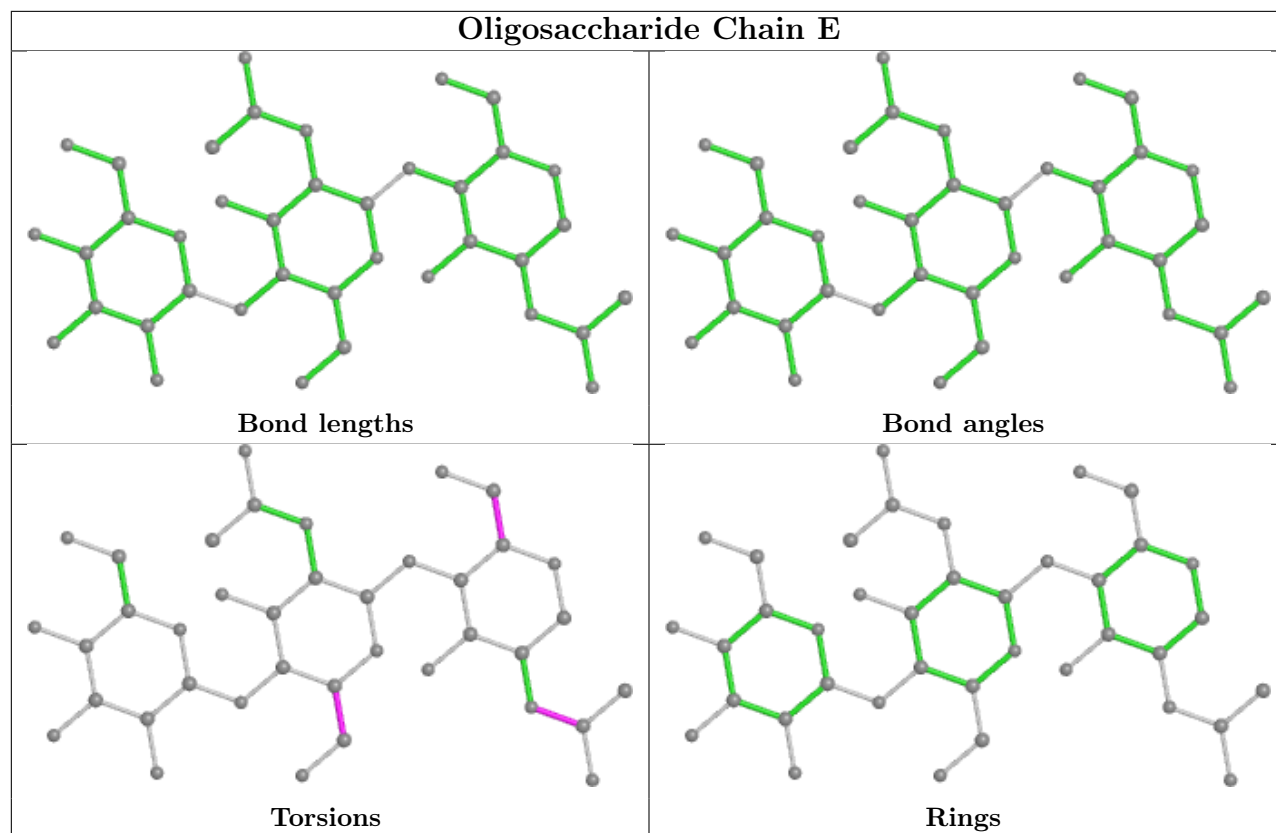


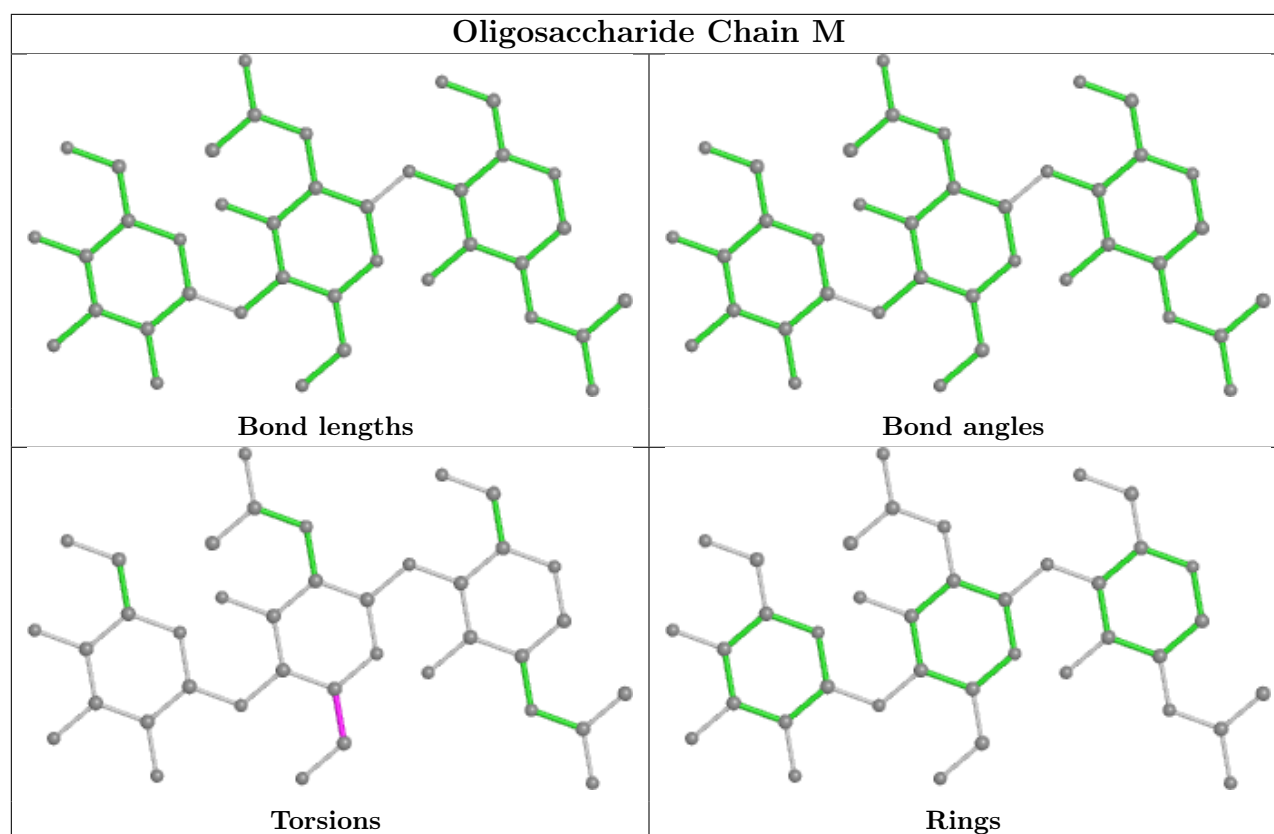












5.6 Ligand geometry [i](#)

24 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	C	1405	1	14,14,15	0.39	0	17,19,21	0.86	1 (5%)
4	NAG	A	1403	1	14,14,15	0.43	0	17,19,21	0.87	1 (5%)
4	NAG	B	1405	1	14,14,15	0.30	0	17,19,21	0.56	0
4	NAG	B	1407	1	14,14,15	0.29	0	17,19,21	0.50	0
4	NAG	C	1406	1	14,14,15	0.25	0	17,19,21	0.46	0
4	NAG	C	1403	1	14,14,15	0.24	0	17,19,21	0.51	0
5	EIC	A	1408	-	19,19,19	0.57	0	19,19,19	0.56	0
4	NAG	B	1402	1	14,14,15	0.24	0	17,19,21	0.47	0
4	NAG	C	1407	1	14,14,15	0.23	0	17,19,21	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	1404	1	14,14,15	0.29	0	17,19,21	0.50	0
4	NAG	A	1401	1	14,14,15	0.28	0	17,19,21	0.43	0
4	NAG	B	1401	1	14,14,15	0.29	0	17,19,21	0.49	0
4	NAG	A	1407	1	14,14,15	0.26	0	17,19,21	0.51	0
5	EIC	C	1401	-	19,19,19	0.56	0	19,19,19	0.57	0
4	NAG	B	1404	1	14,14,15	0.30	0	17,19,21	0.52	0
5	EIC	B	1408	-	19,19,19	0.57	0	19,19,19	0.61	0
4	NAG	C	1402	1	14,14,15	0.37	0	17,19,21	0.49	0
4	NAG	A	1406	1	14,14,15	0.28	0	17,19,21	0.48	0
4	NAG	B	1406	1	14,14,15	0.40	0	17,19,21	0.37	0
4	NAG	A	1405	1	14,14,15	0.26	0	17,19,21	0.54	0
4	NAG	C	1408	1	14,14,15	0.32	0	17,19,21	0.47	0
4	NAG	B	1403	1	14,14,15	0.31	0	17,19,21	0.54	0
4	NAG	C	1404	1	14,14,15	0.26	0	17,19,21	0.50	0
4	NAG	A	1402	1	14,14,15	0.29	0	17,19,21	0.48	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	C	1405	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1403	1	-	1/6/23/26	0/1/1/1
4	NAG	B	1405	1	-	4/6/23/26	0/1/1/1
4	NAG	B	1407	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1406	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1403	1	-	2/6/23/26	0/1/1/1
5	EIC	A	1408	-	-	2/17/17/17	-
4	NAG	B	1402	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1407	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1404	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1401	1	-	1/6/23/26	0/1/1/1
4	NAG	B	1401	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1407	1	-	2/6/23/26	0/1/1/1
5	EIC	C	1401	-	-	4/17/17/17	-
4	NAG	B	1404	1	-	2/6/23/26	0/1/1/1
5	EIC	B	1408	-	-	7/17/17/17	-
4	NAG	C	1402	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1406	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	1406	1	-	1/6/23/26	0/1/1/1
4	NAG	A	1405	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1408	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1403	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1404	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1402	1	-	2/6/23/26	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(°)
4	C	1405	NAG	C2-N2-C7	2.47	126.42	122.90
4	A	1403	NAG	C2-N2-C7	2.38	126.29	122.90

There are no chirality outliers.

All (46) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1405	NAG	O5-C5-C6-O6
4	C	1403	NAG	O5-C5-C6-O6
4	C	1404	NAG	O5-C5-C6-O6
4	B	1405	NAG	C4-C5-C6-O6
4	A	1402	NAG	O5-C5-C6-O6
4	C	1404	NAG	C4-C5-C6-O6
4	A	1404	NAG	O5-C5-C6-O6
4	A	1407	NAG	O5-C5-C6-O6
4	C	1406	NAG	O5-C5-C6-O6
4	C	1403	NAG	C4-C5-C6-O6
4	B	1405	NAG	C8-C7-N2-C2
4	B	1405	NAG	O7-C7-N2-C2
4	A	1405	NAG	O5-C5-C6-O6
4	B	1401	NAG	O5-C5-C6-O6
4	C	1402	NAG	O5-C5-C6-O6
4	A	1402	NAG	C4-C5-C6-O6
4	A	1404	NAG	C4-C5-C6-O6
4	C	1406	NAG	C4-C5-C6-O6
4	A	1406	NAG	O5-C5-C6-O6
4	C	1402	NAG	C4-C5-C6-O6
4	A	1407	NAG	C4-C5-C6-O6
4	B	1401	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	A	1406	NAG	C4-C5-C6-O6
4	B	1404	NAG	O5-C5-C6-O6
5	C	1401	EIC	C2-C3-C4-C5
4	C	1408	NAG	O5-C5-C6-O6
4	B	1406	NAG	O5-C5-C6-O6
4	C	1408	NAG	C4-C5-C6-O6
5	C	1401	EIC	C13-C14-C15-C16
5	C	1401	EIC	C6-C7-C8-C9
4	A	1401	NAG	O5-C5-C6-O6
5	B	1408	EIC	C10-C11-C12-C13
5	C	1401	EIC	C10-C11-C12-C13
5	B	1408	EIC	C1-C2-C3-C4
5	A	1408	EIC	C3-C4-C5-C6
5	A	1408	EIC	C7-C8-C9-C10
5	B	1408	EIC	C15-C16-C17-C18
4	A	1403	NAG	C3-C2-N2-C7
4	C	1405	NAG	C3-C2-N2-C7
5	B	1408	EIC	C6-C7-C8-C9
5	B	1408	EIC	O1-C1-C2-C3
5	B	1408	EIC	O2-C1-C2-C3
5	B	1408	EIC	C2-C3-C4-C5
4	B	1404	NAG	C4-C5-C6-O6
4	A	1405	NAG	C4-C5-C6-O6
4	C	1405	NAG	C4-C5-C6-O6

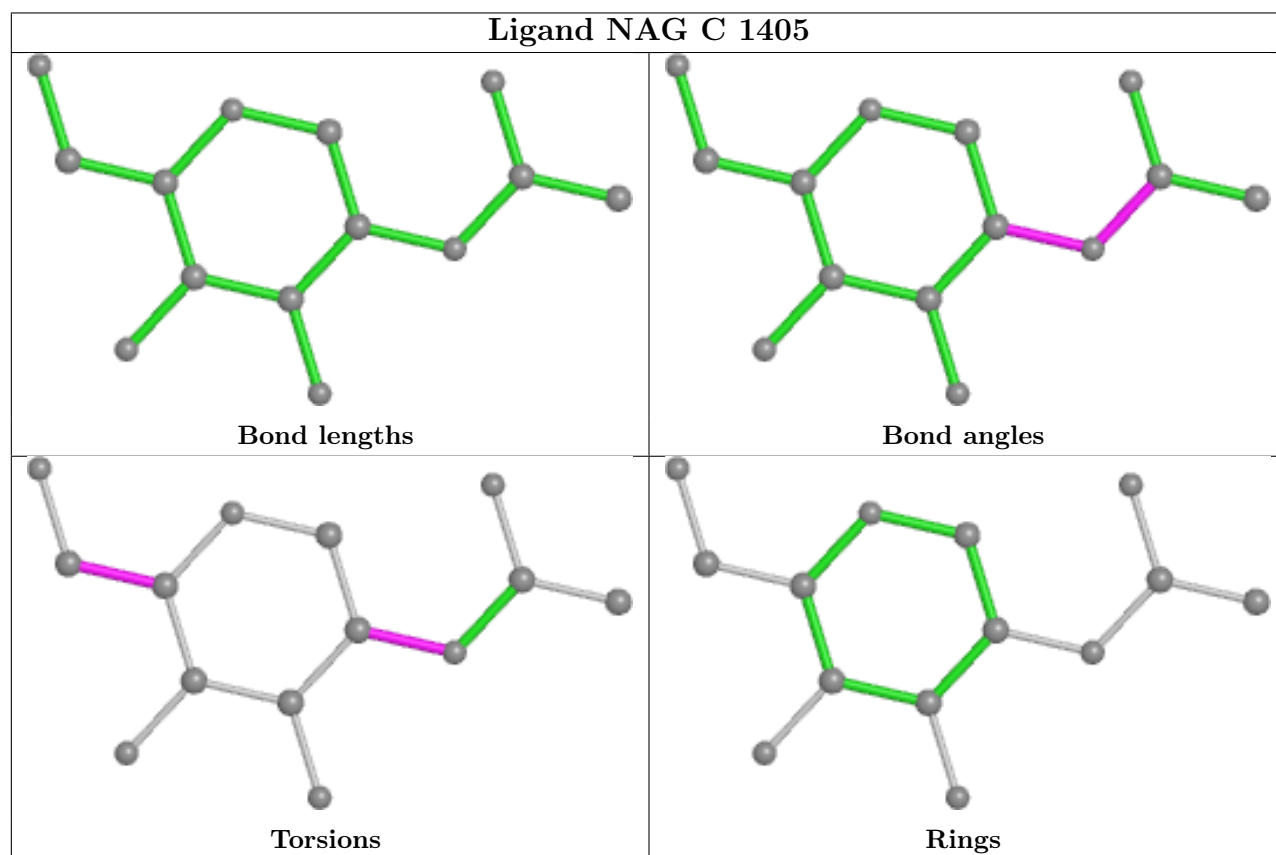
There are no ring outliers.

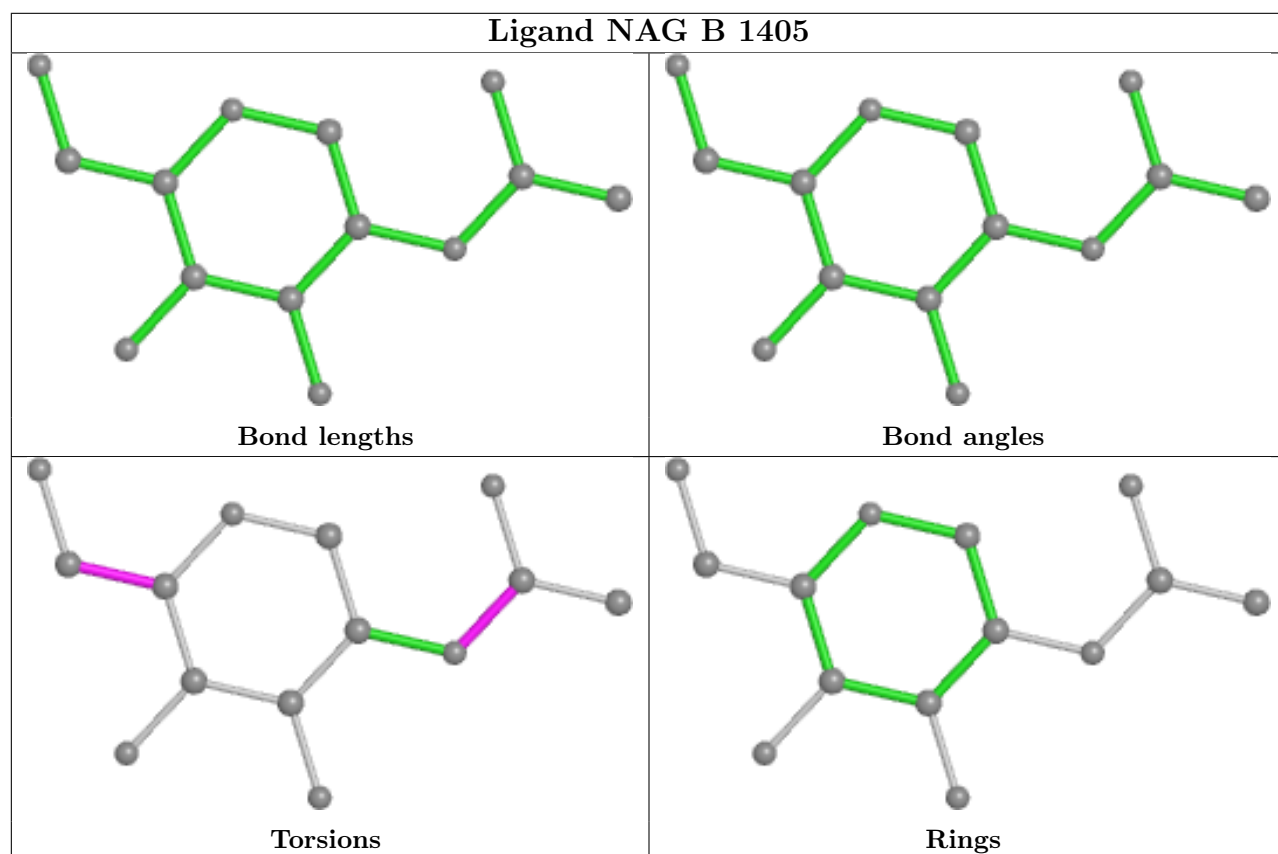
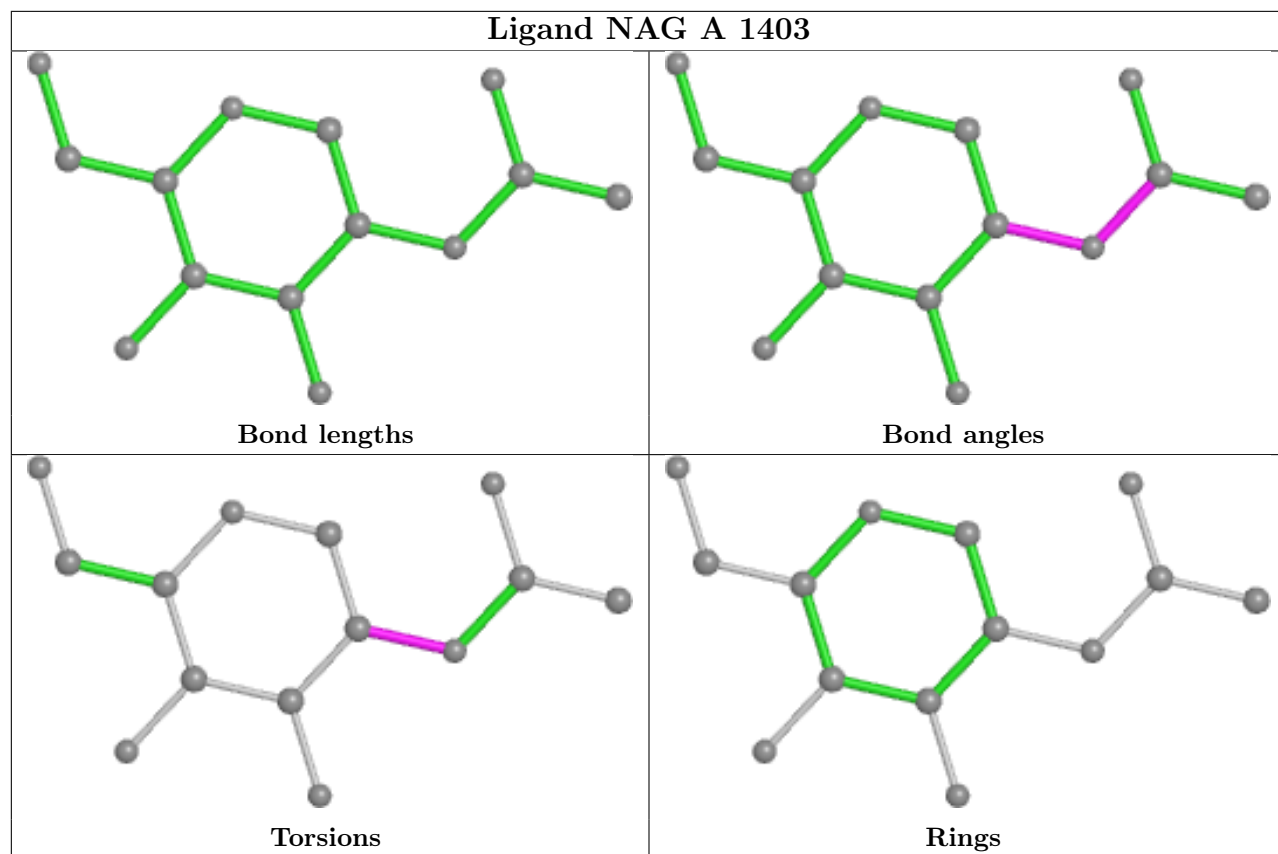
6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1408	EIC	2	0
5	C	1401	EIC	3	0
4	B	1404	NAG	1	0
5	B	1408	EIC	2	0
4	A	1405	NAG	1	0
4	B	1403	NAG	1	0

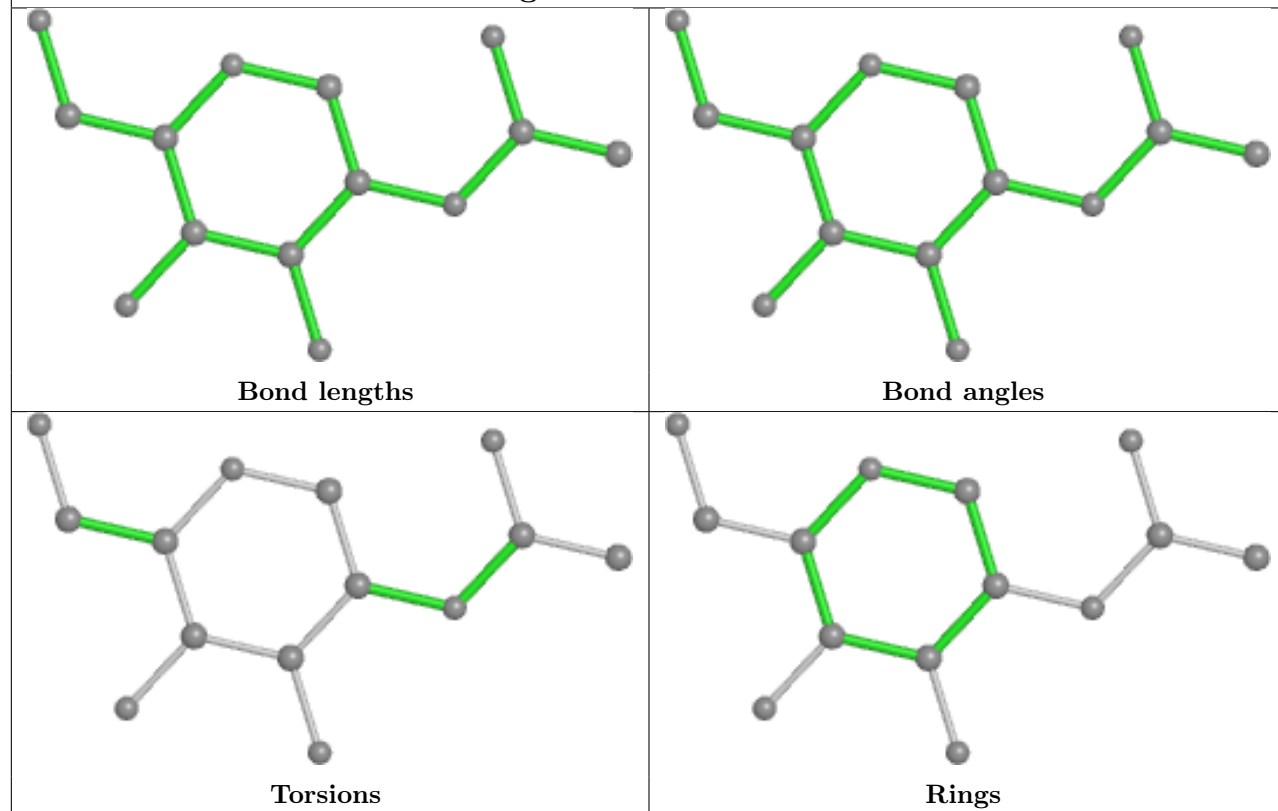
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

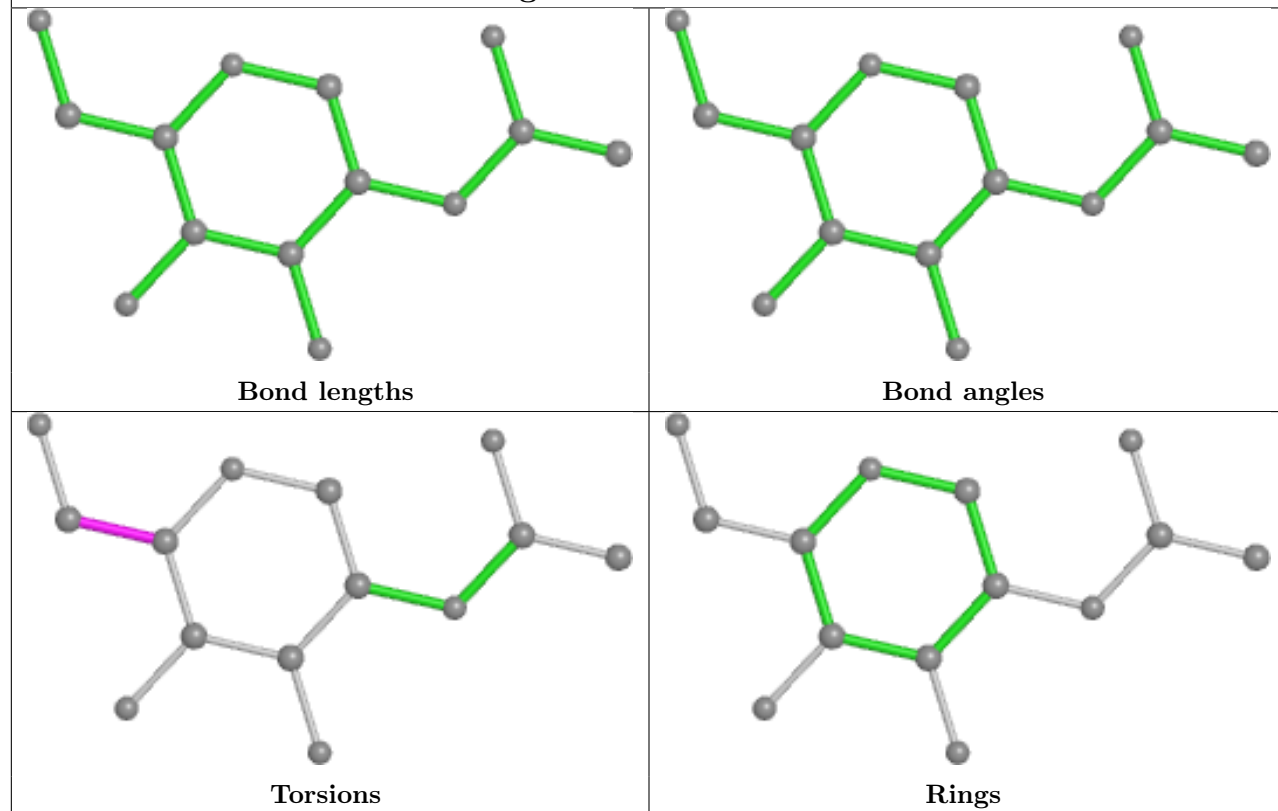


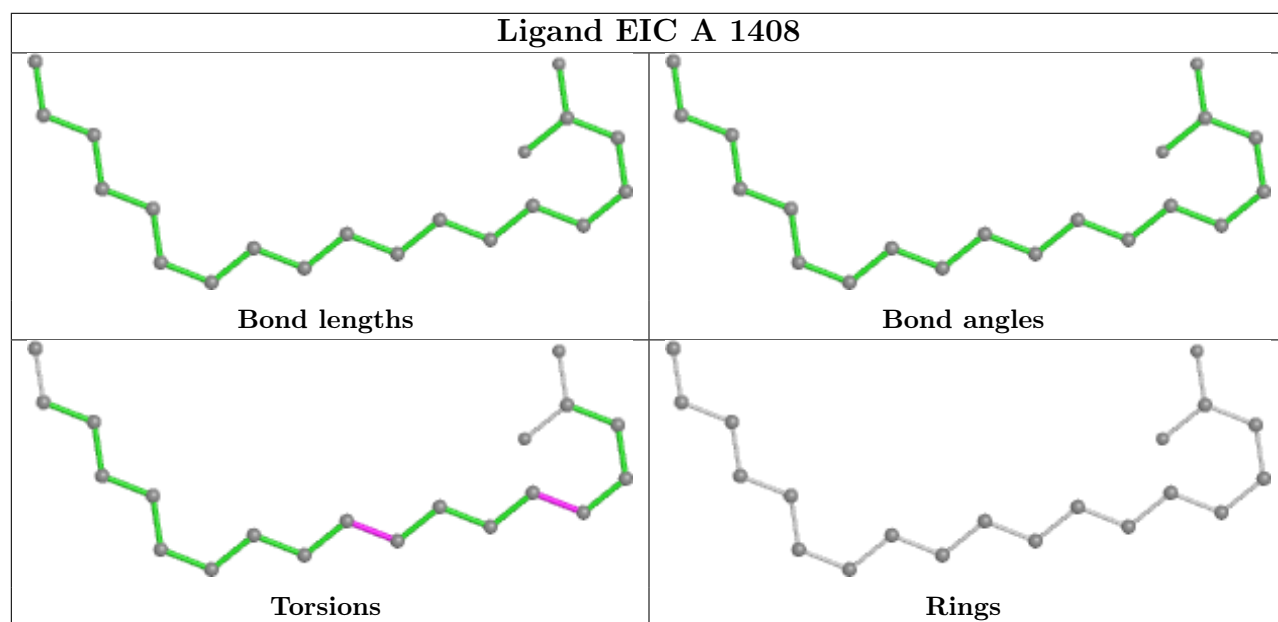
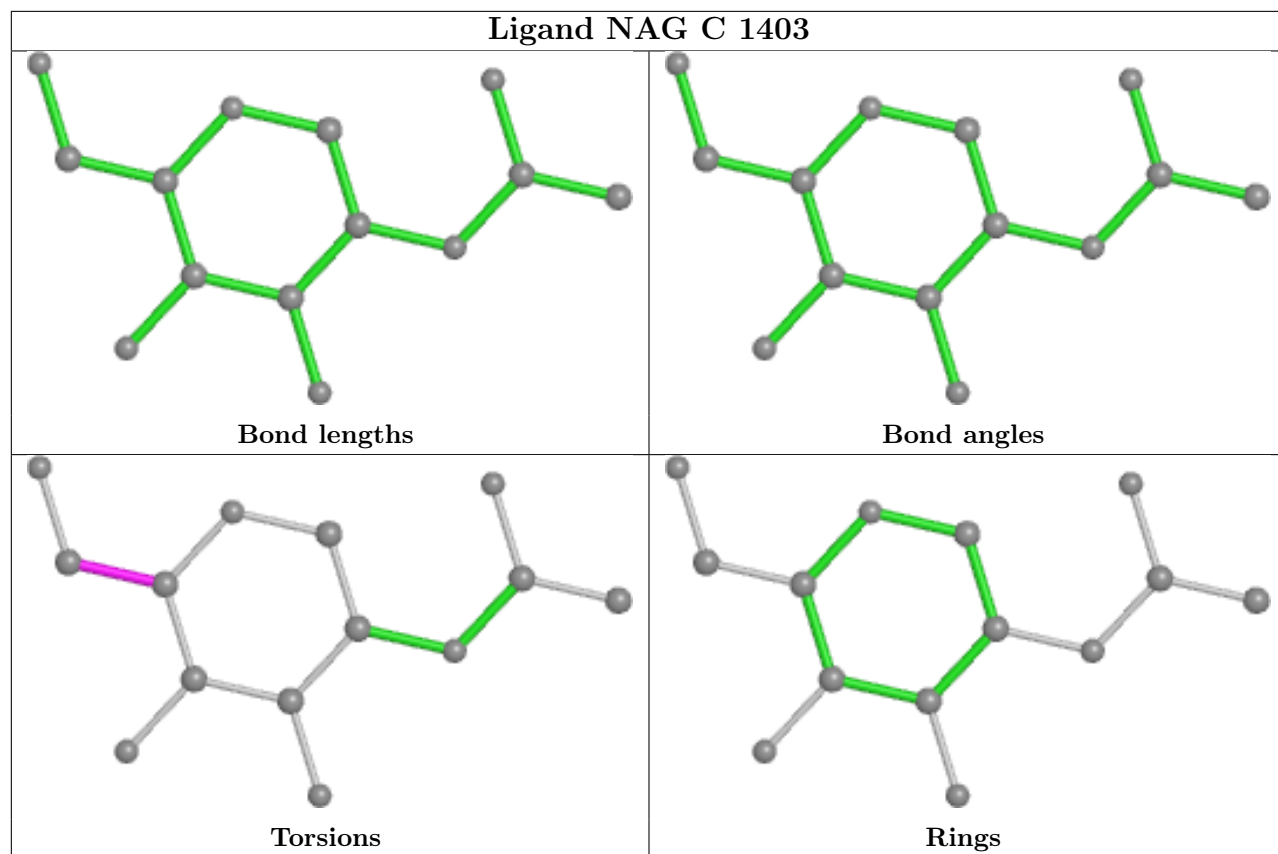


Ligand NAG B 1407

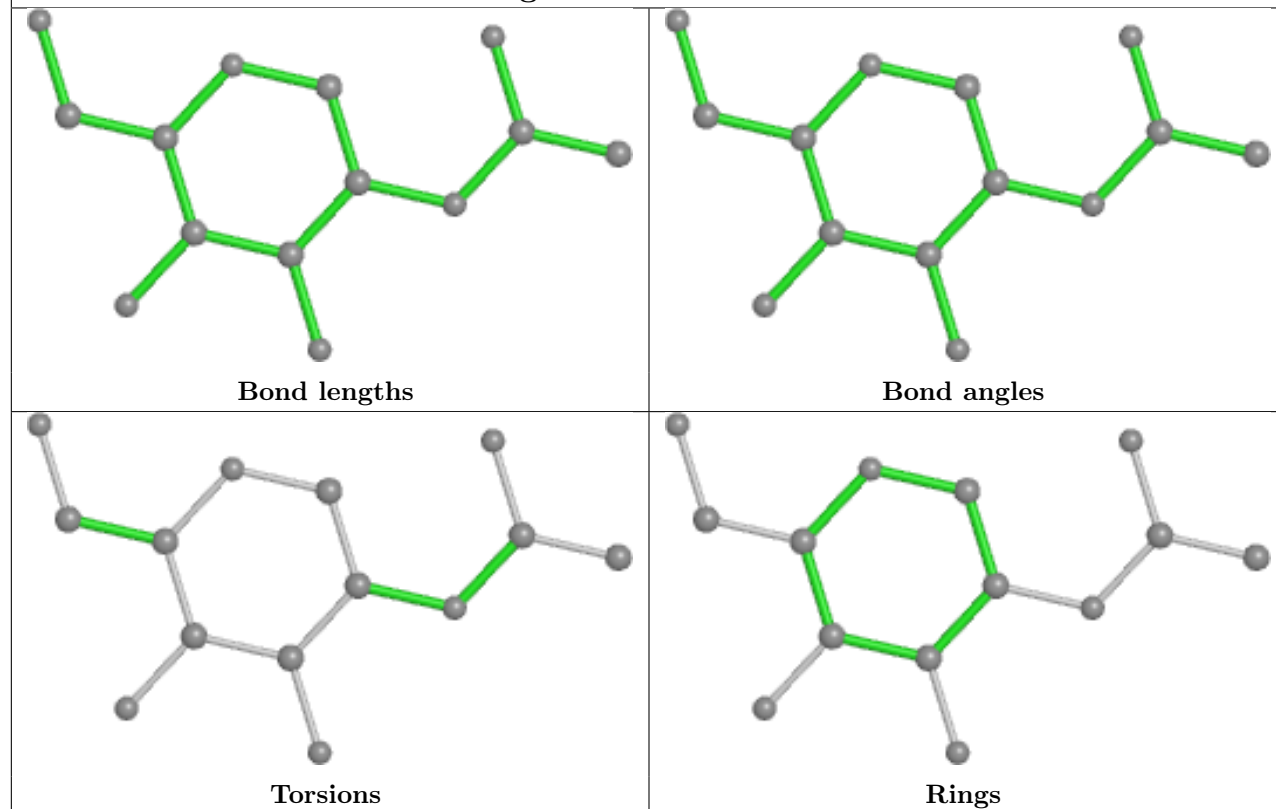


Ligand NAG C 1406

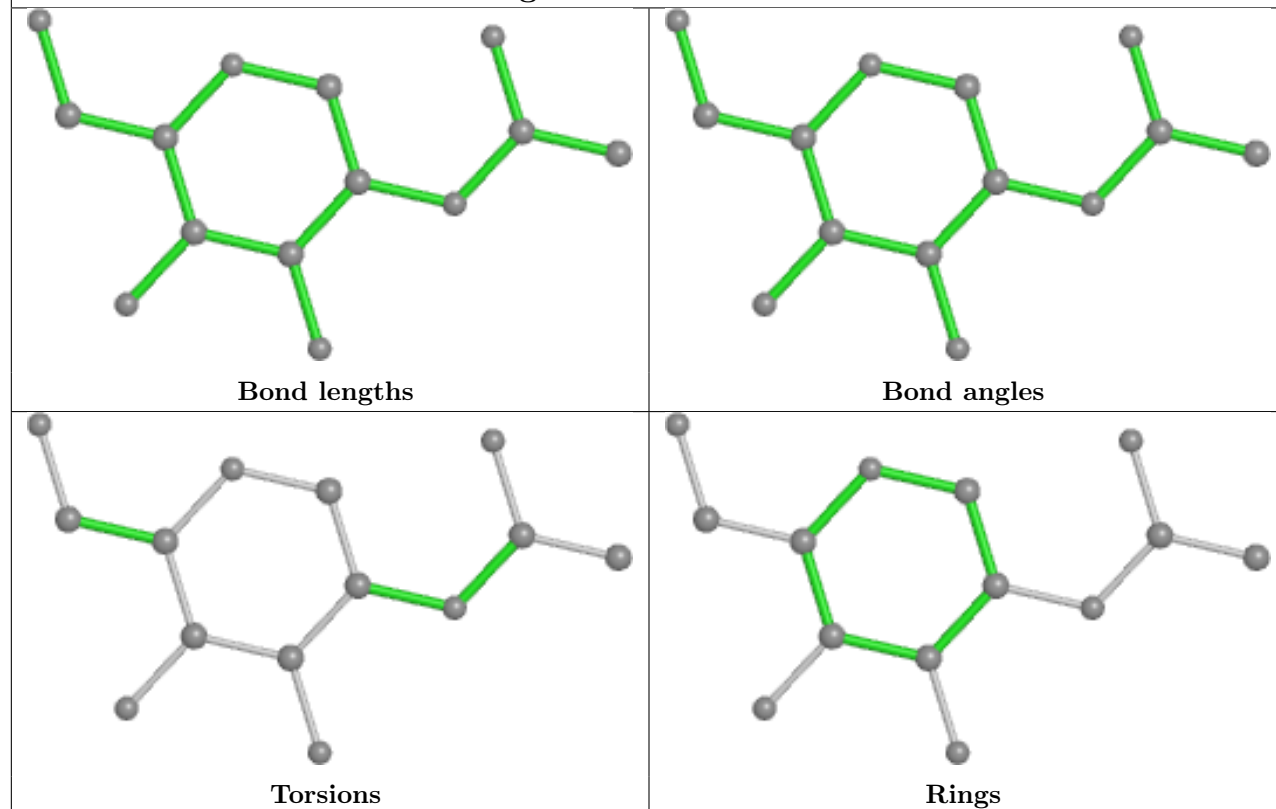


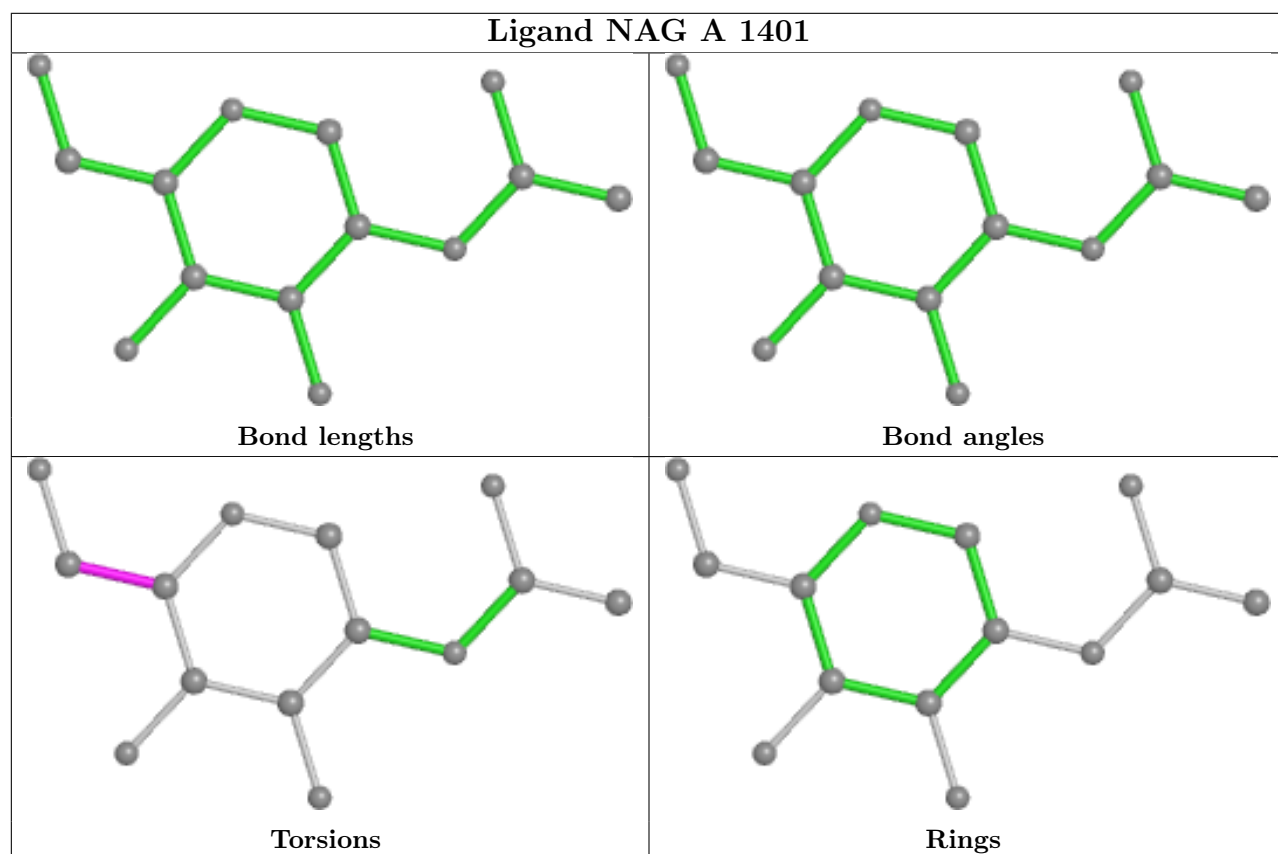
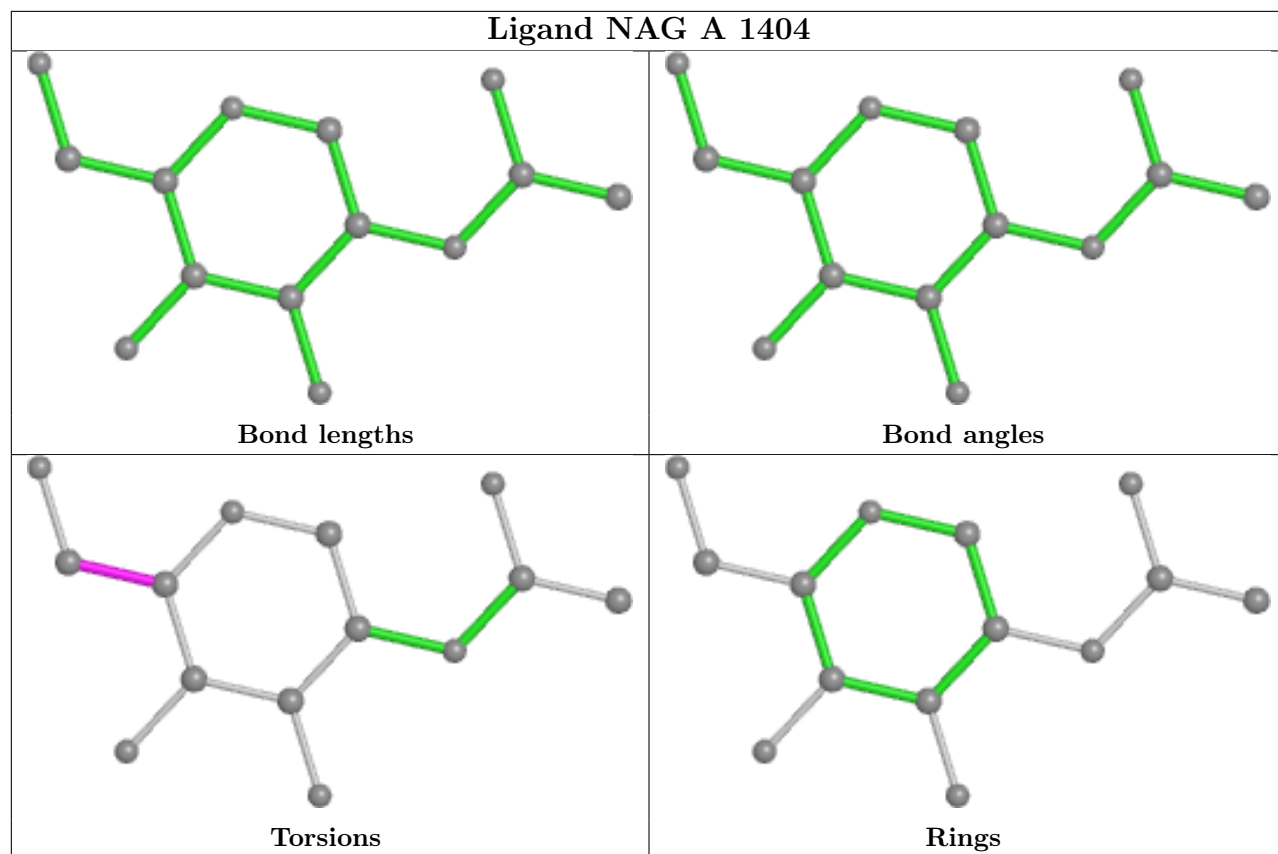


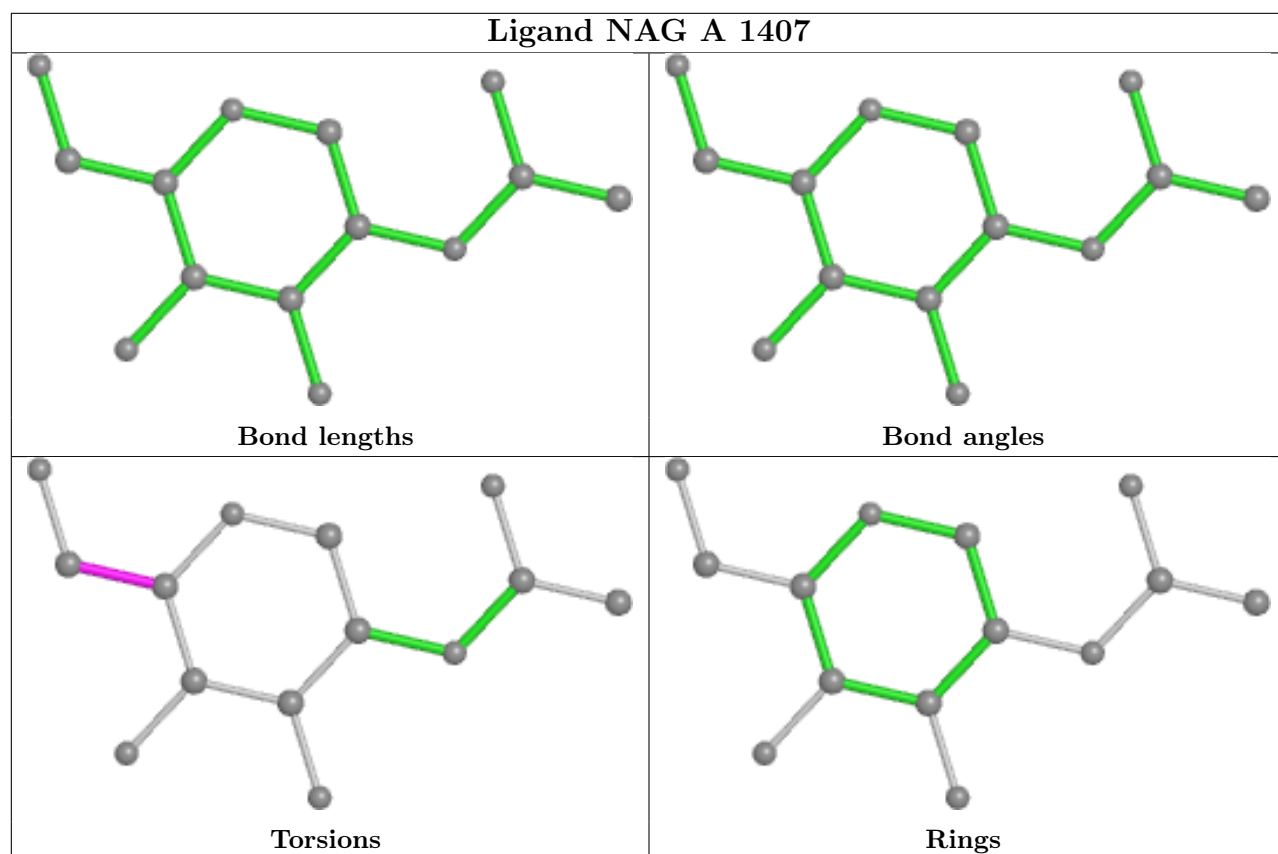
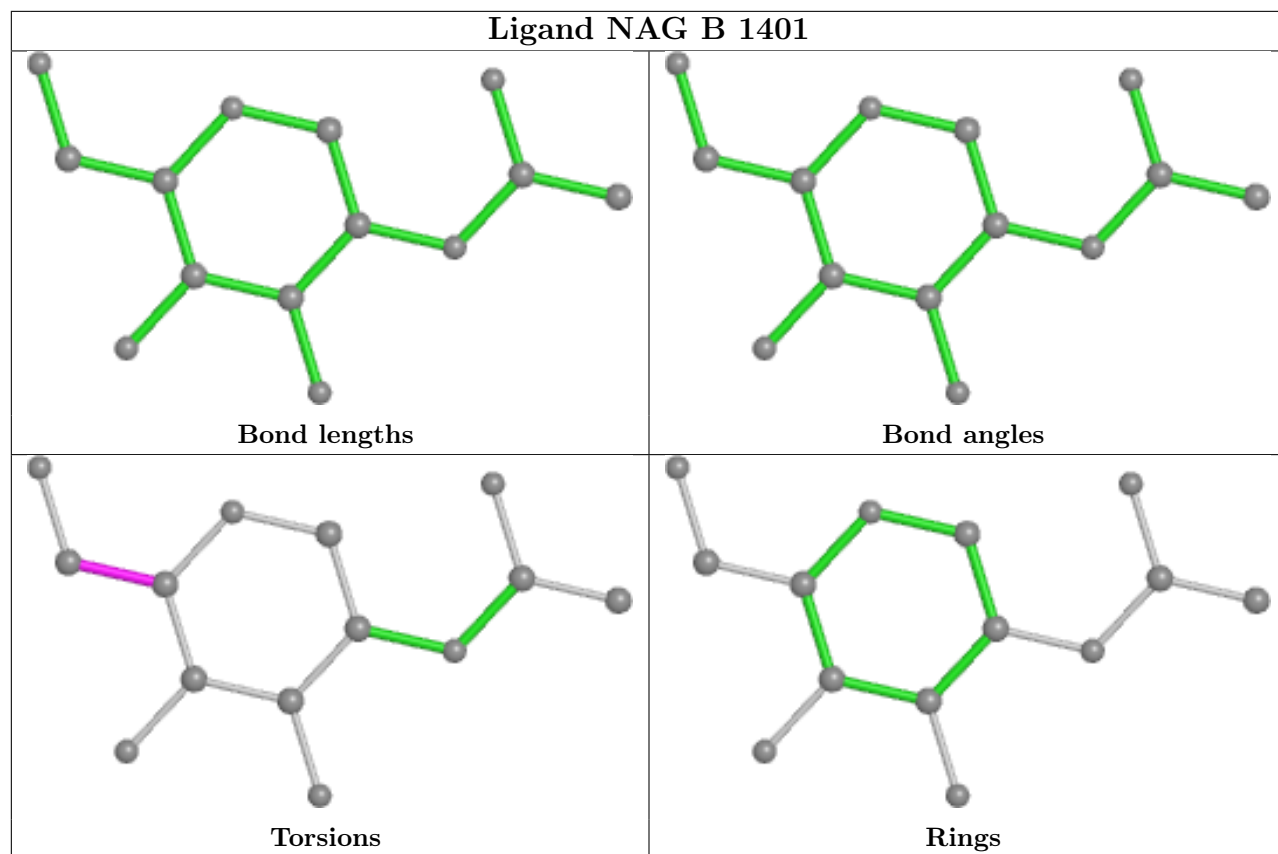
Ligand NAG B 1402

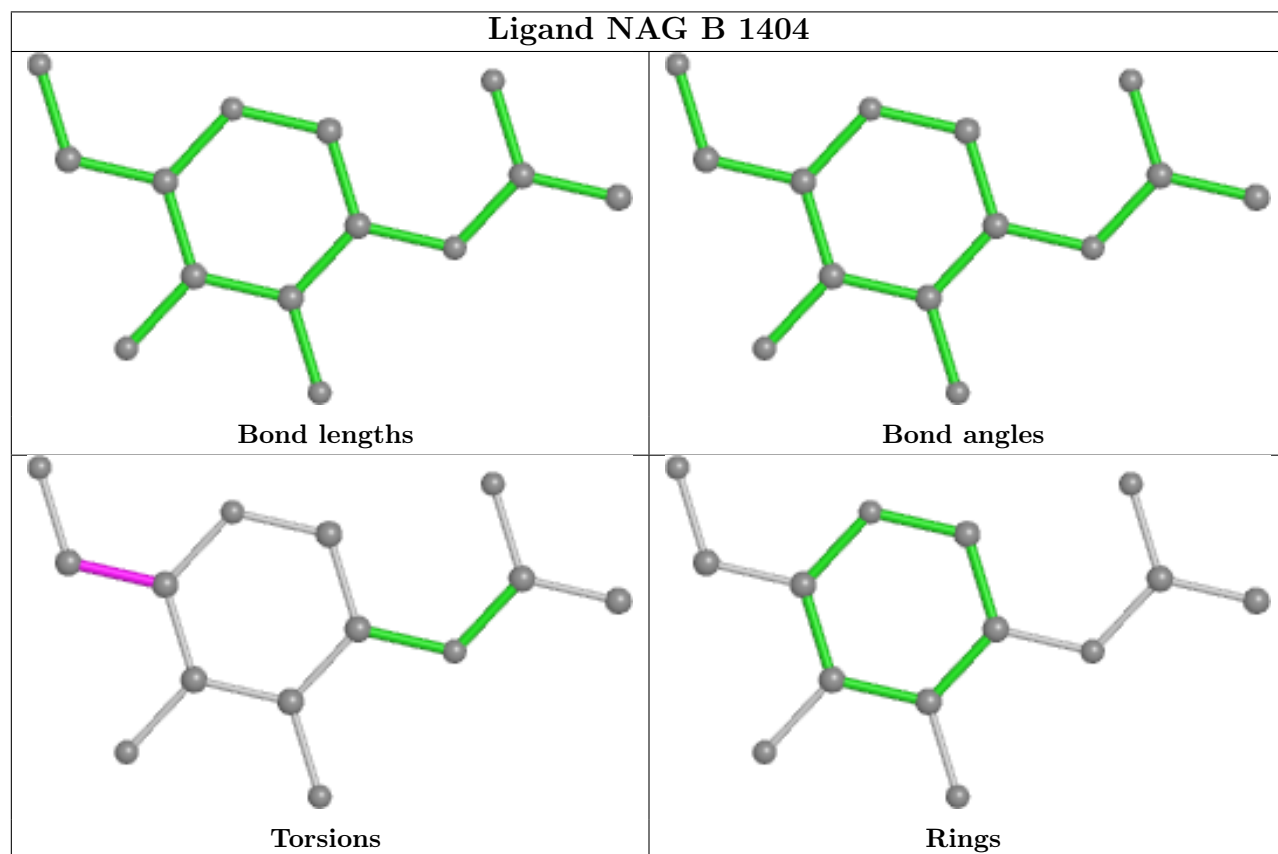
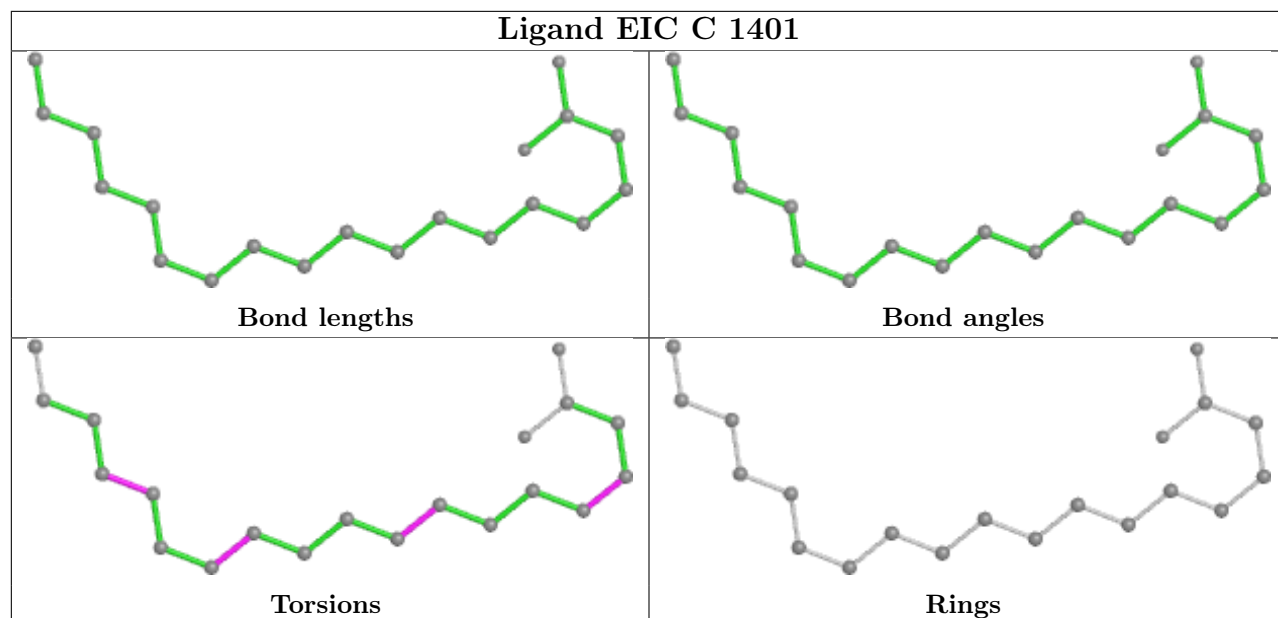


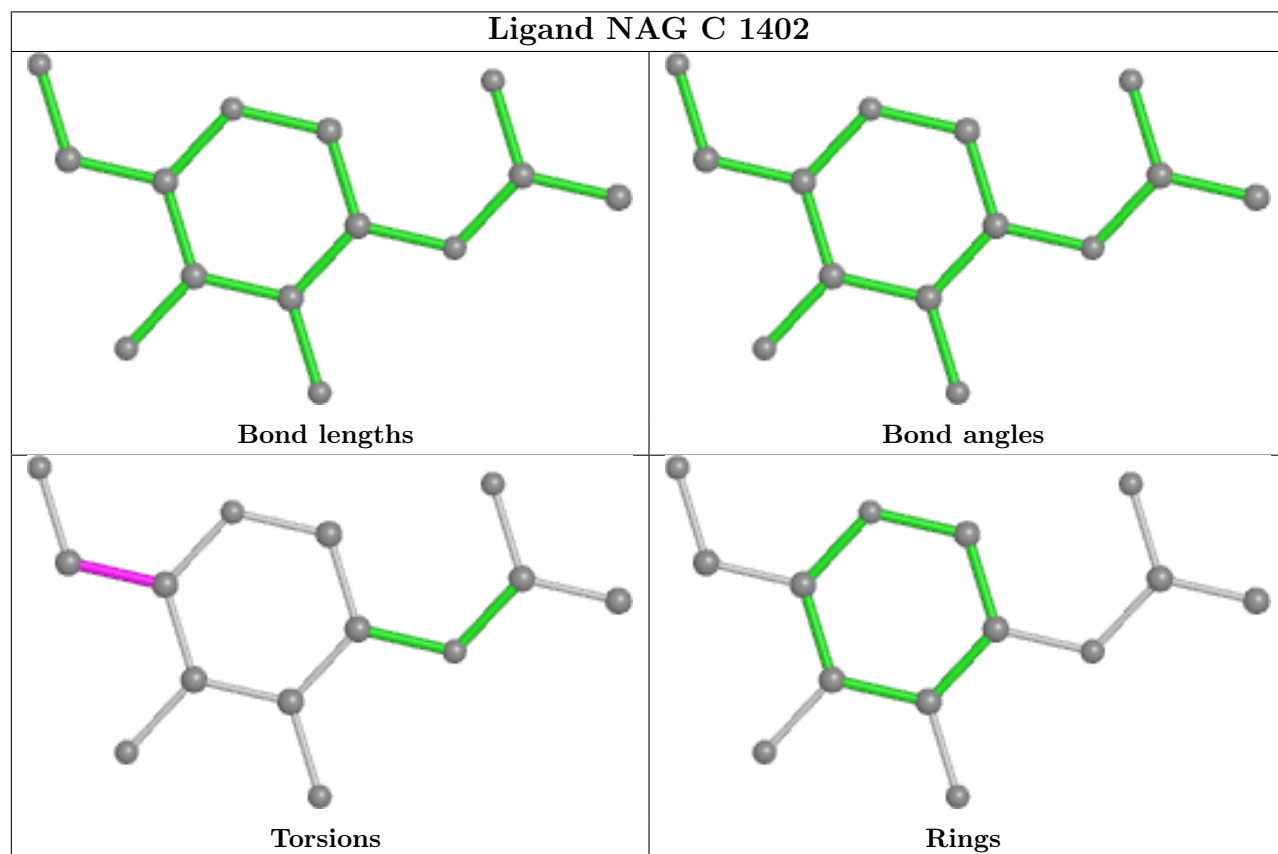
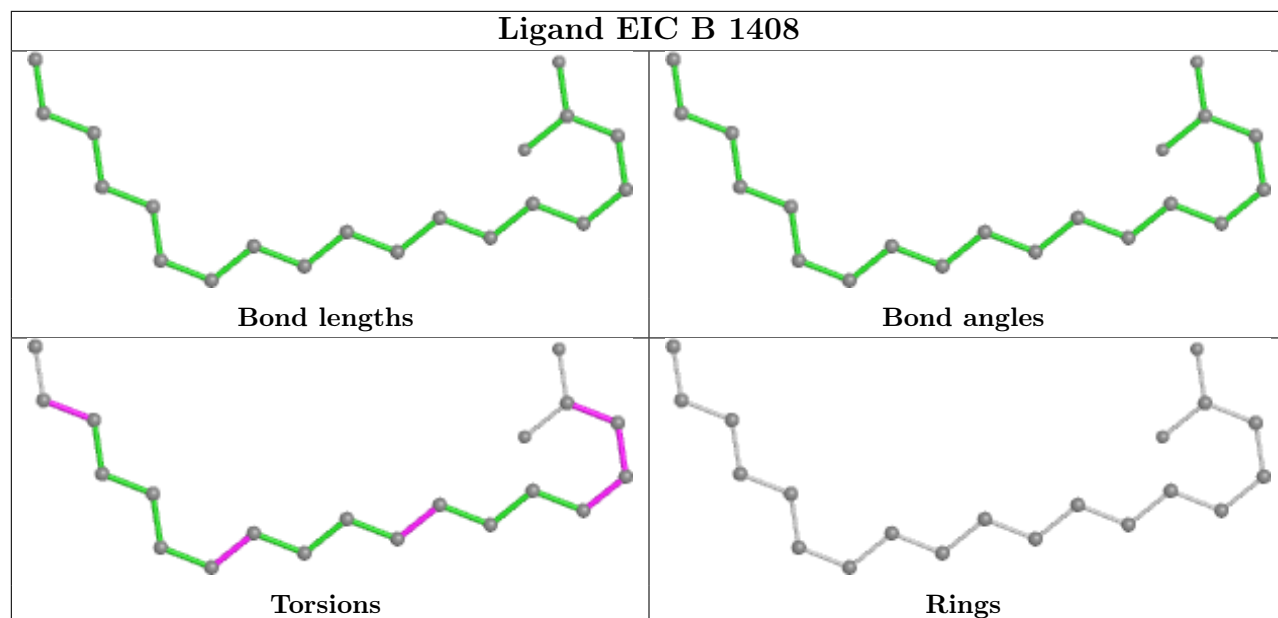
Ligand NAG C 1407



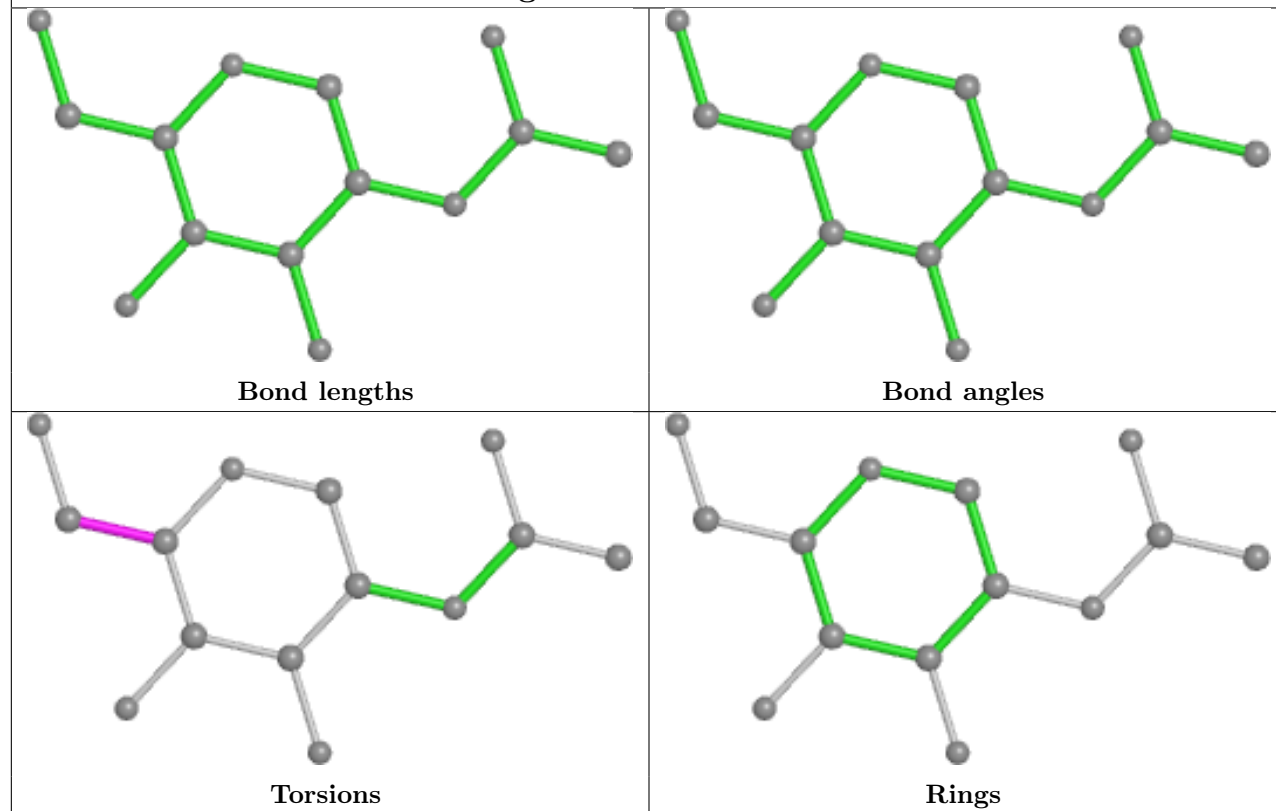




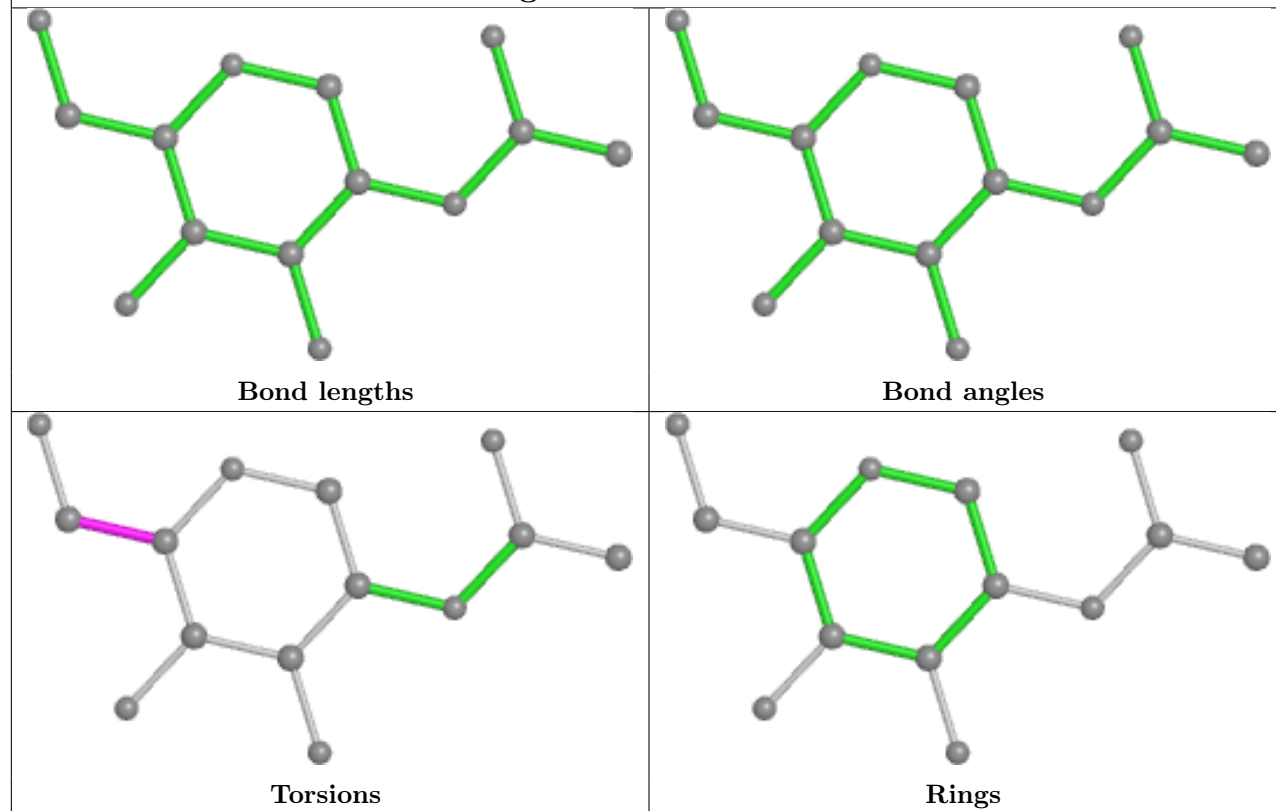




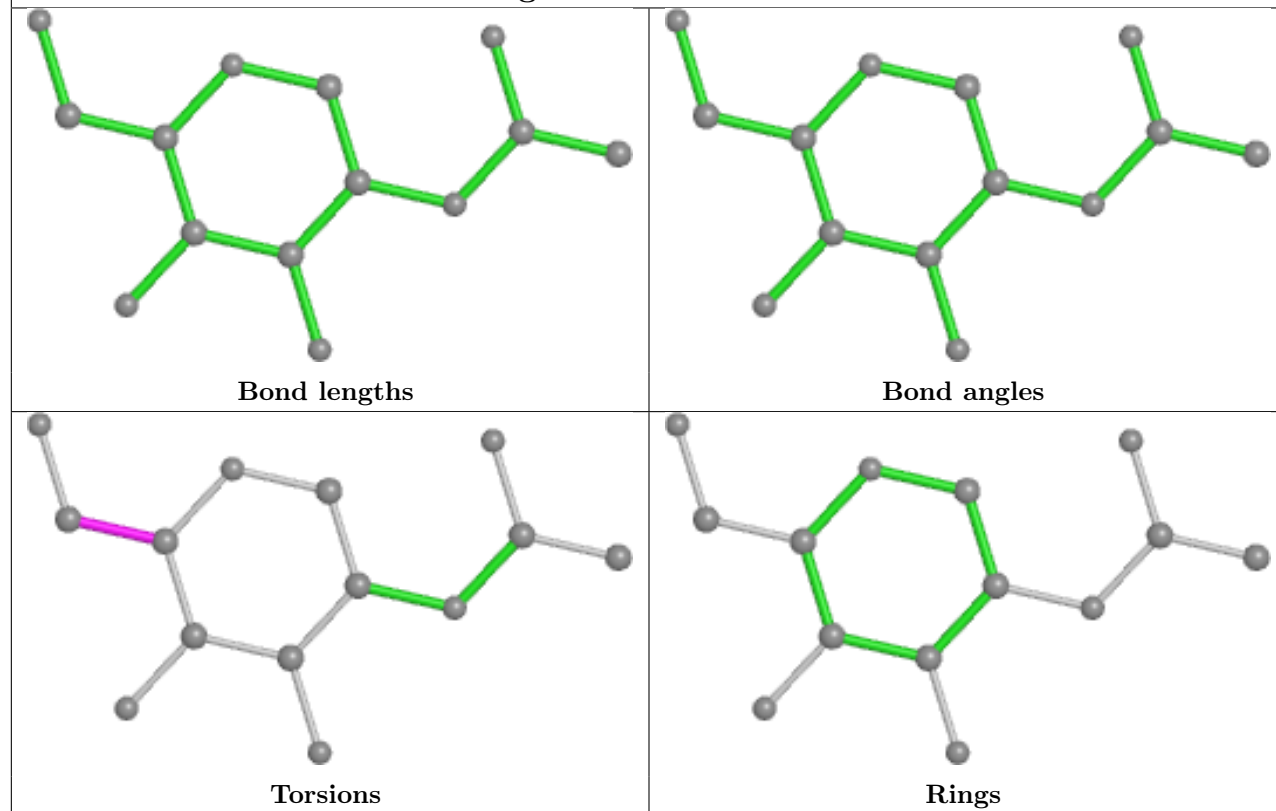
Ligand NAG A 1406



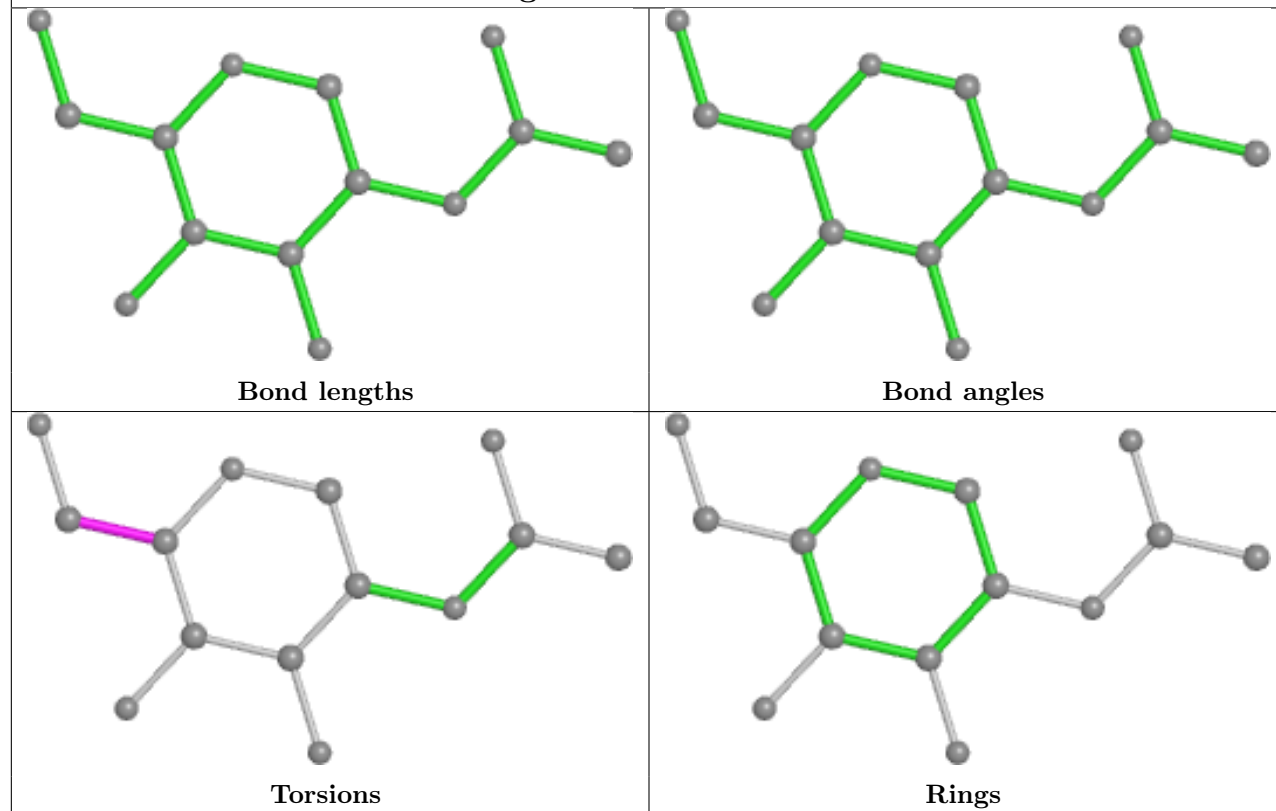
Ligand NAG B 1406

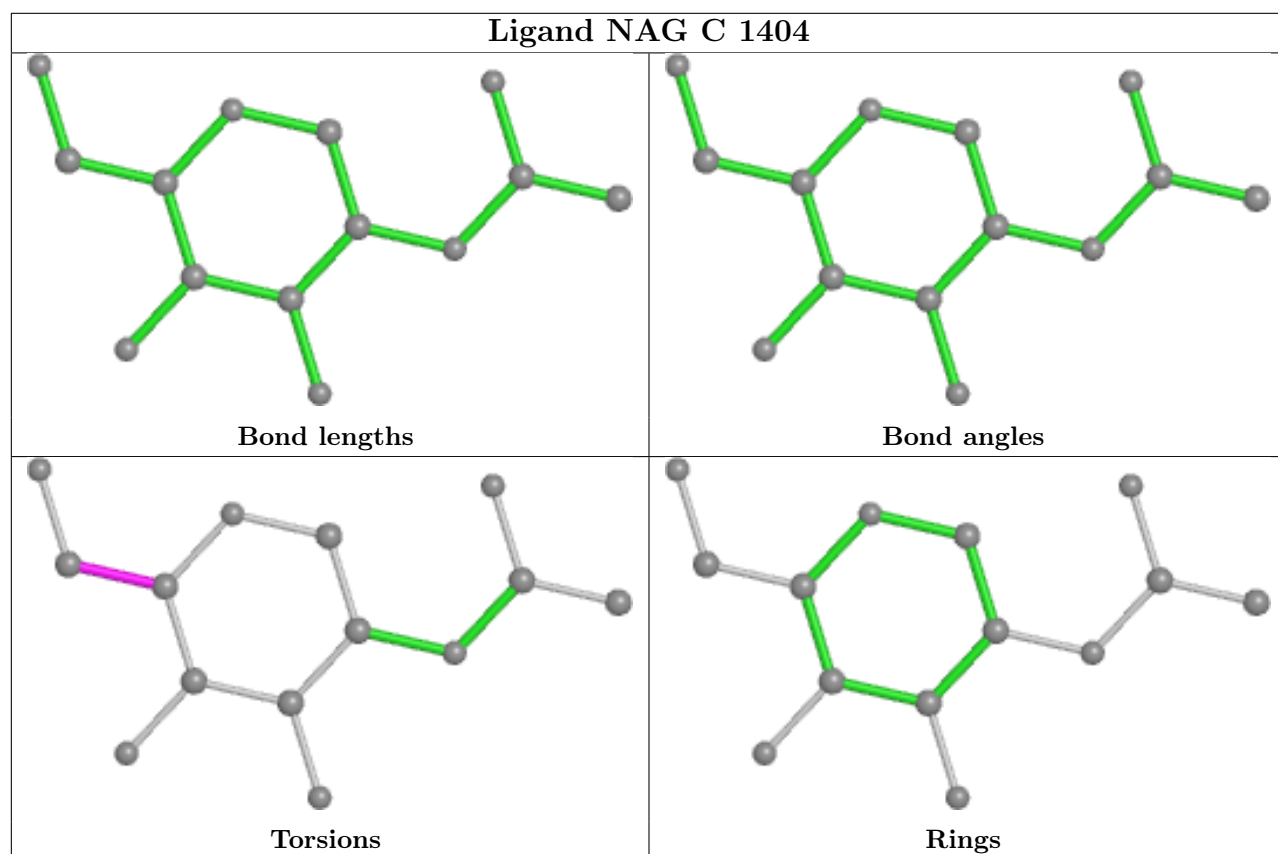
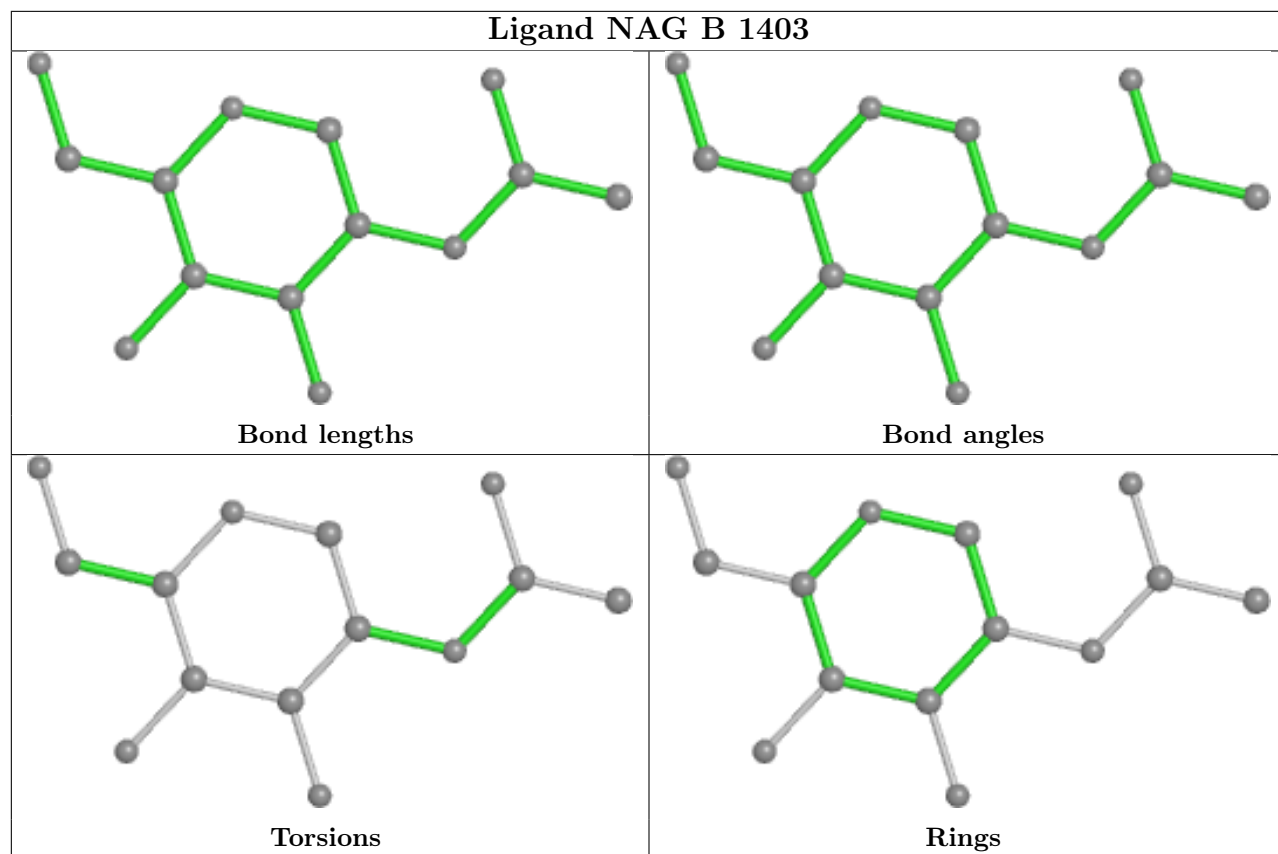


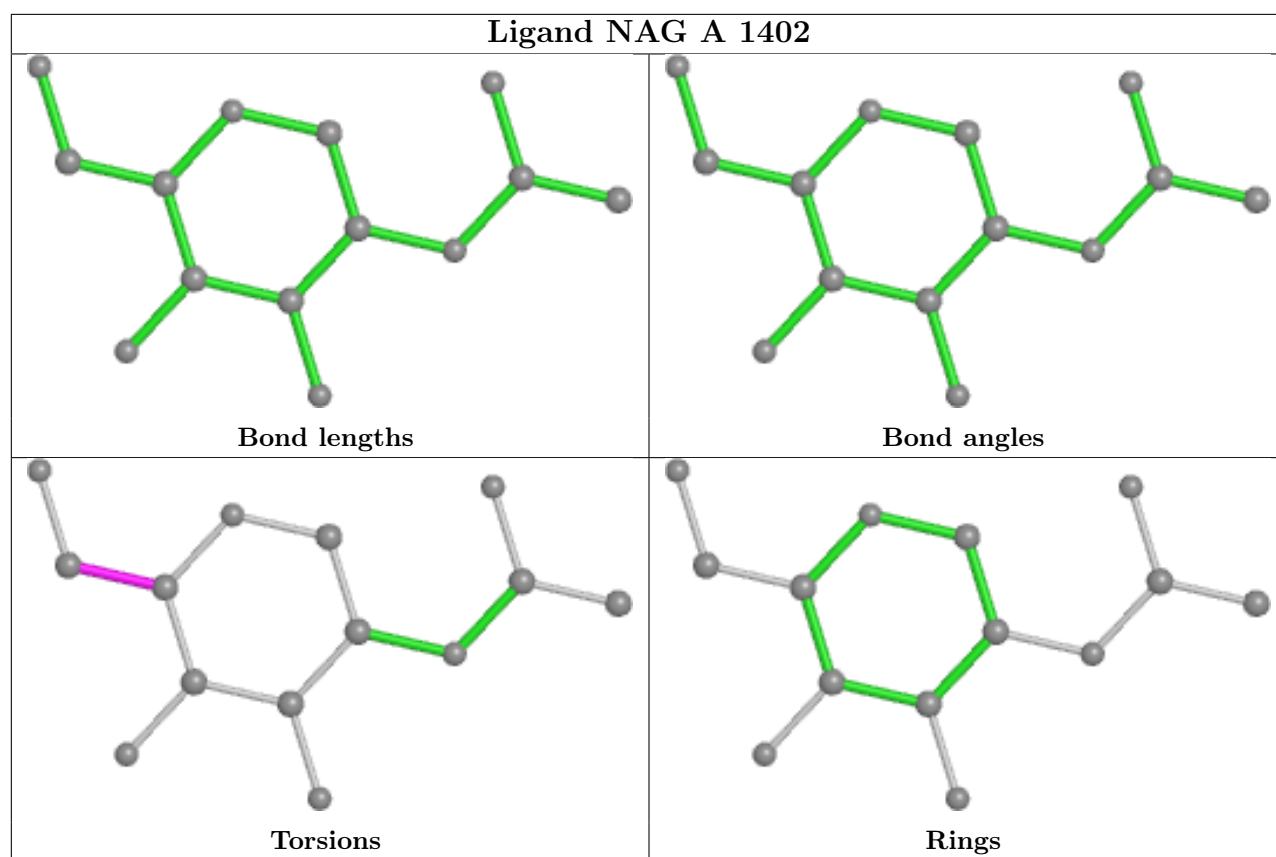
Ligand NAG A 1405



Ligand NAG C 1408







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.

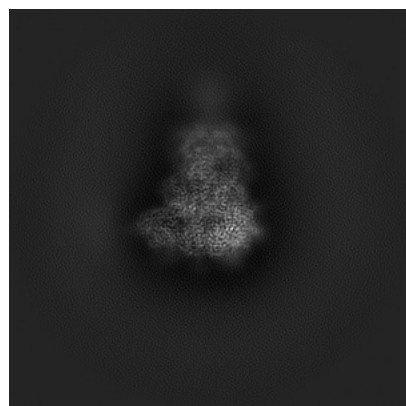
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-61609. These allow visual inspection of the internal detail of the map and identification of artifacts.

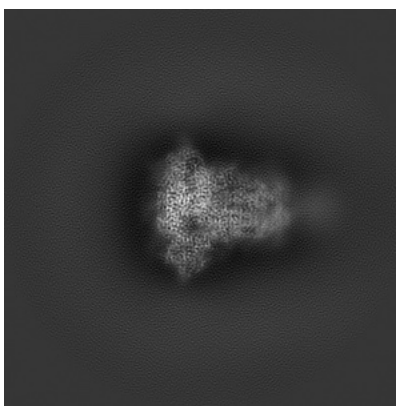
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

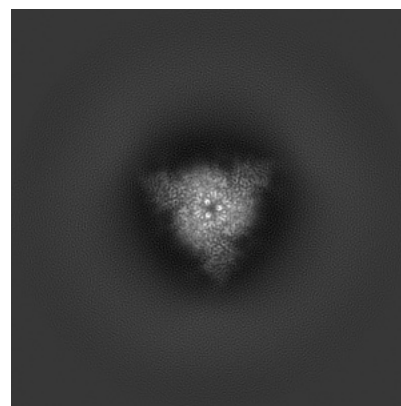
6.1.1 Primary map



X

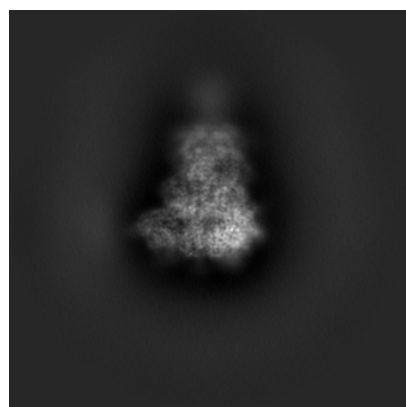


Y

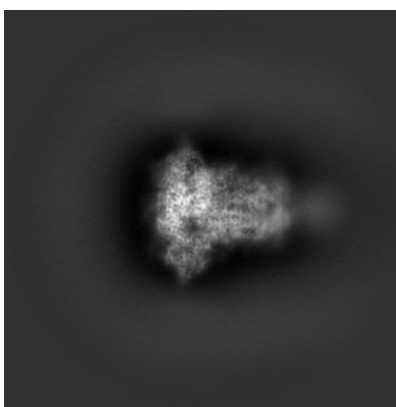


Z

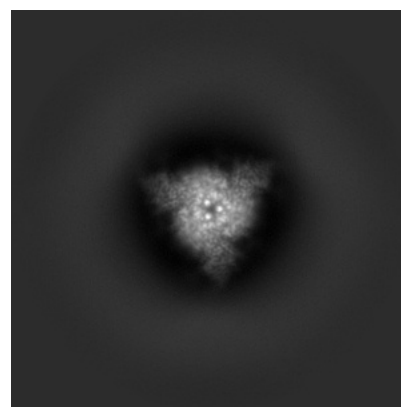
6.1.2 Raw map



X



Y

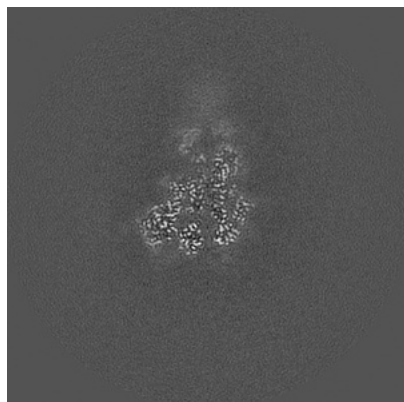


Z

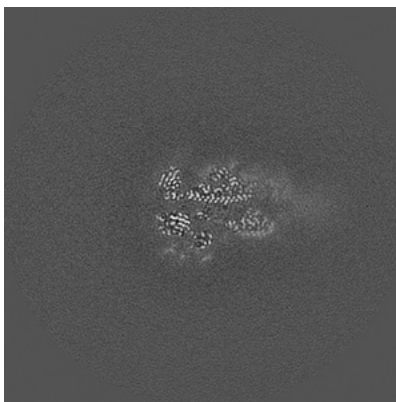
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

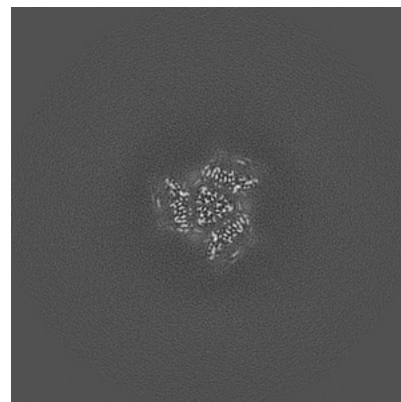
6.2.1 Primary map



X Index: 160

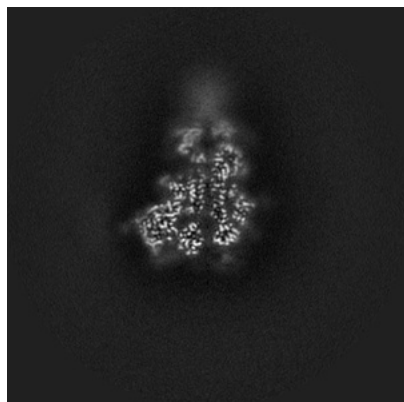


Y Index: 160

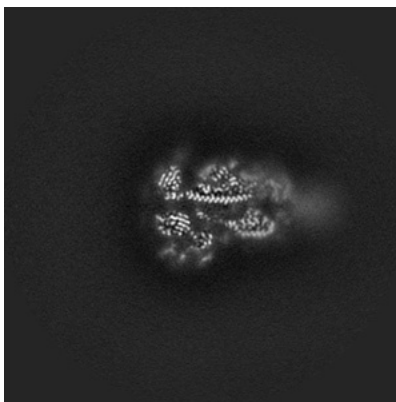


Z Index: 160

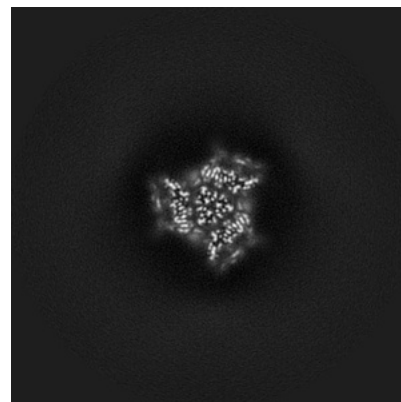
6.2.2 Raw map



X Index: 160



Y Index: 160

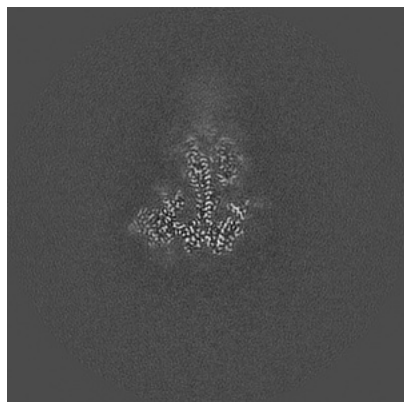


Z Index: 160

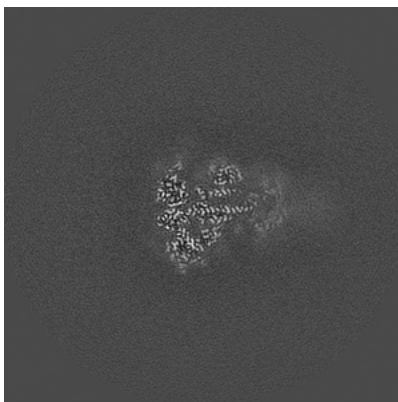
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

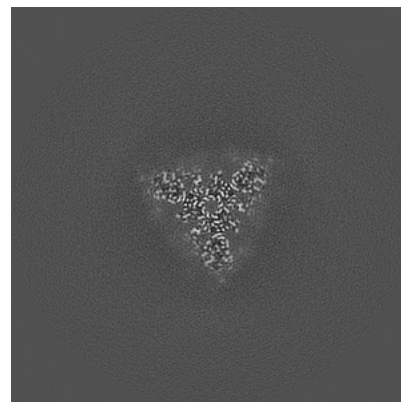
6.3.1 Primary map



X Index: 167

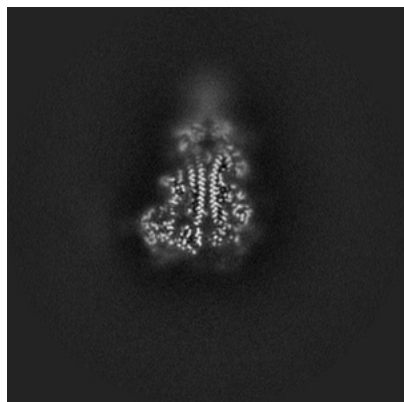


Y Index: 166

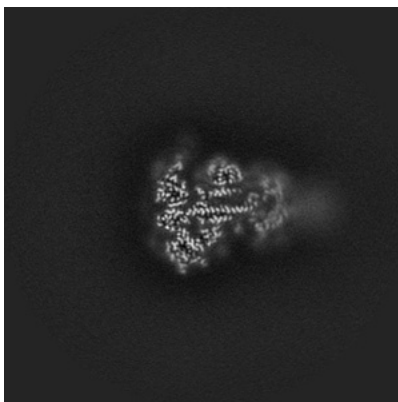


Z Index: 139

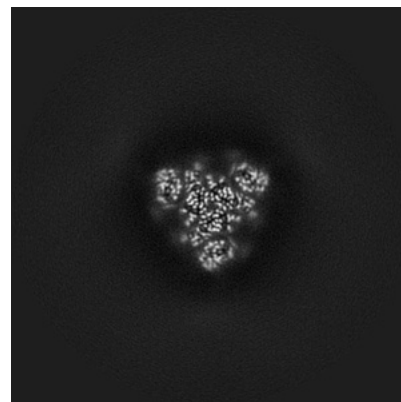
6.3.2 Raw map



X Index: 157



Y Index: 166

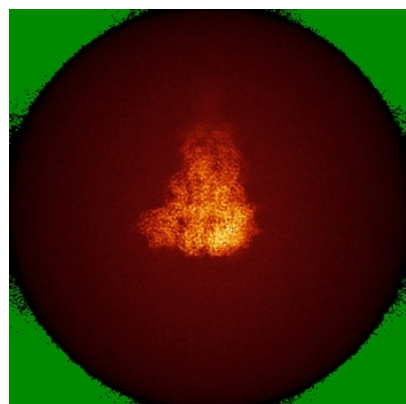


Z Index: 134

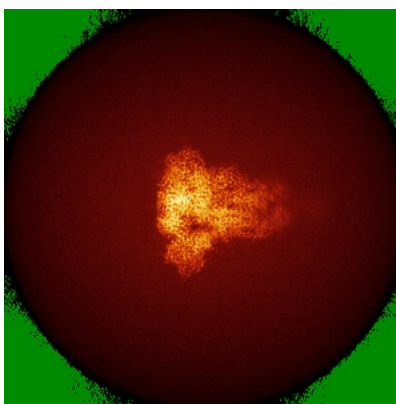
The images above show the largest variance slices of the map in three orthogonal directions.

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

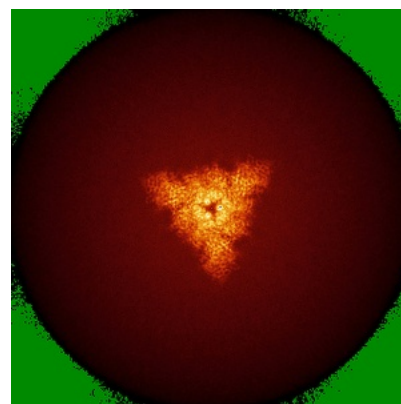
6.4.1 Primary map



X



Y

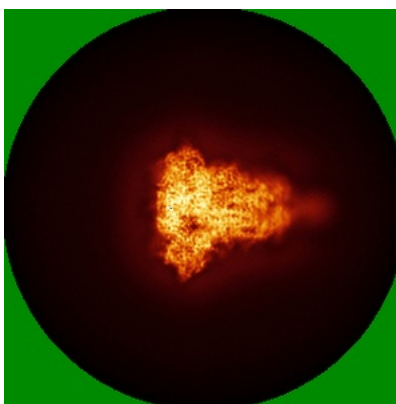


Z

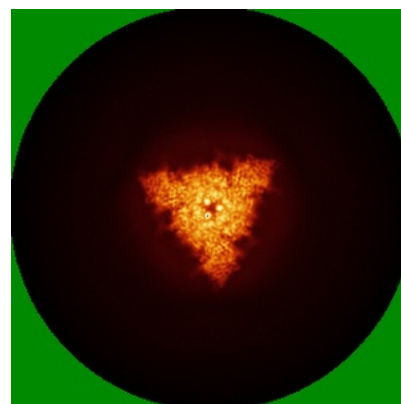
6.4.2 Raw map



X



Y

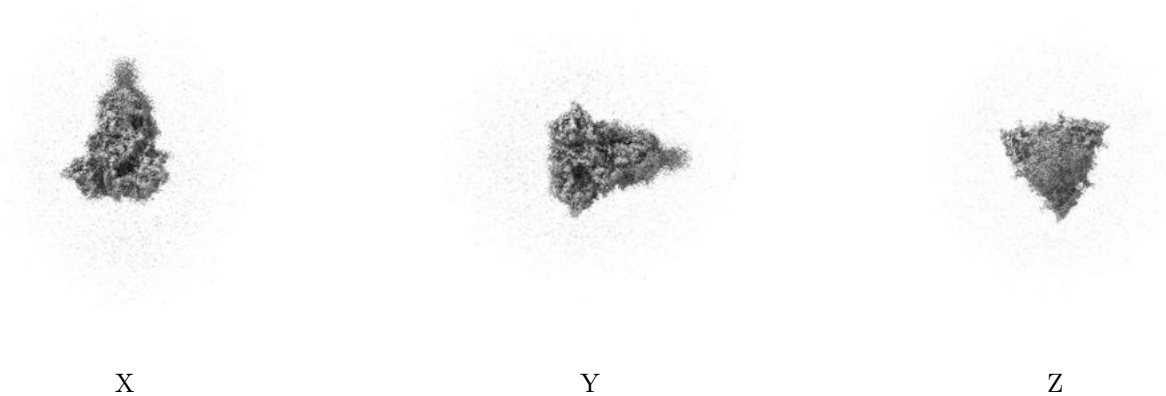


Z

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.

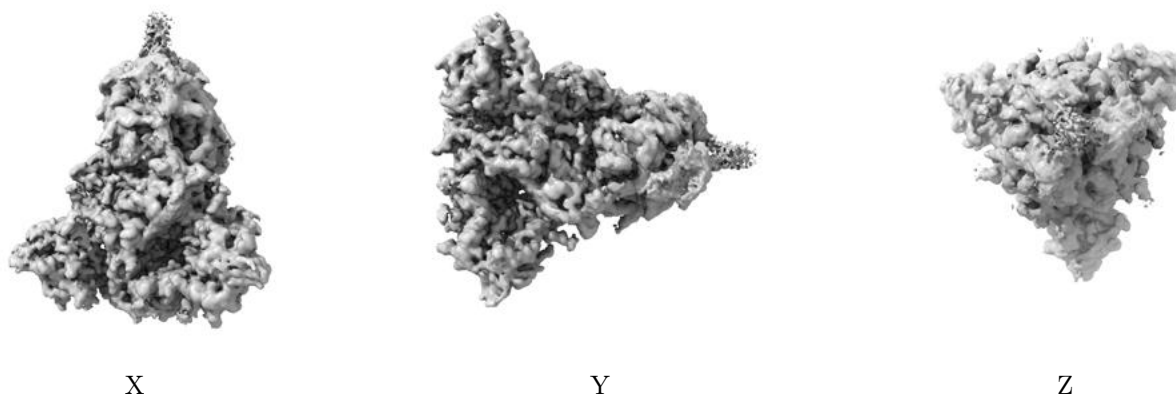
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.022. 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.

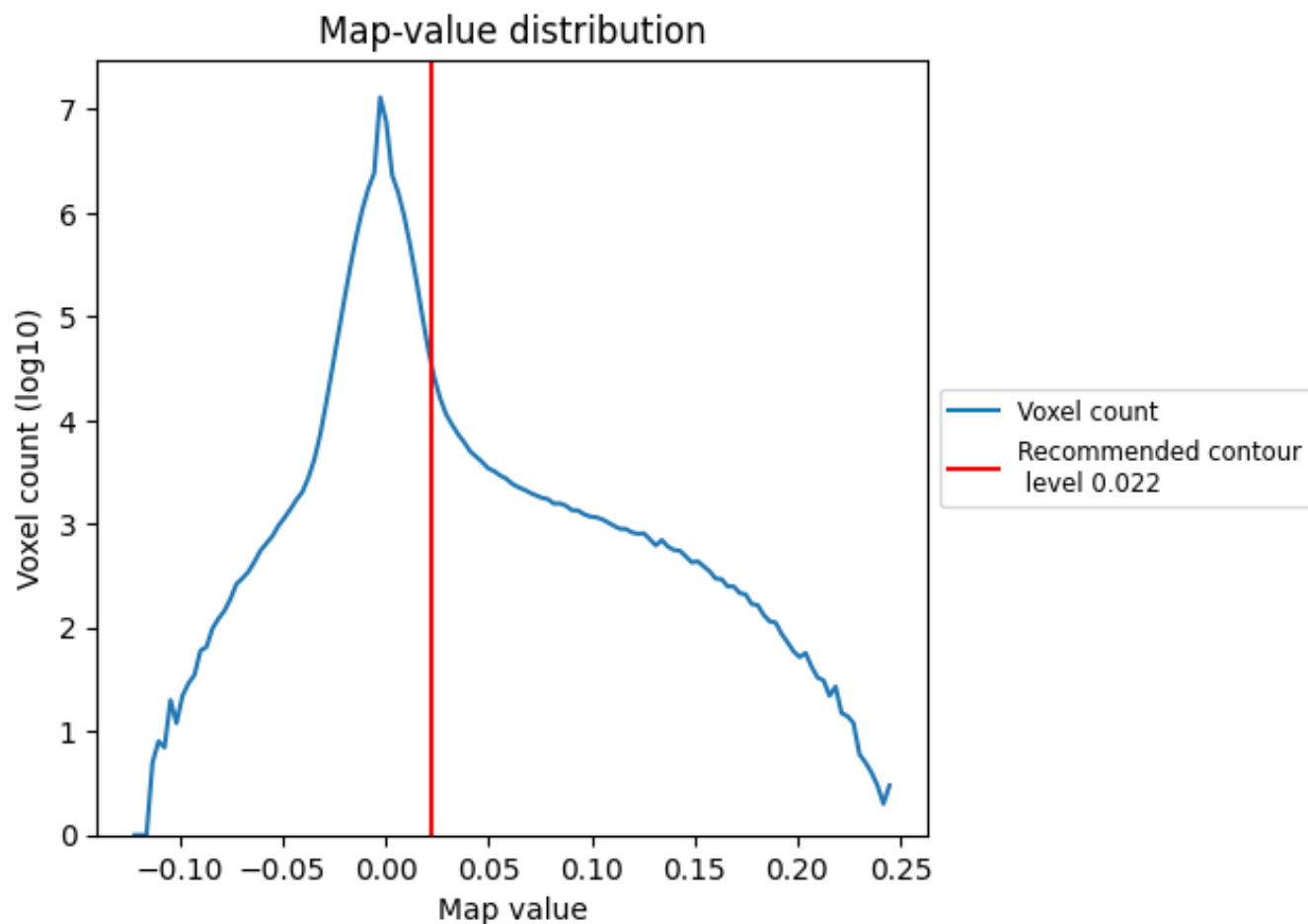
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

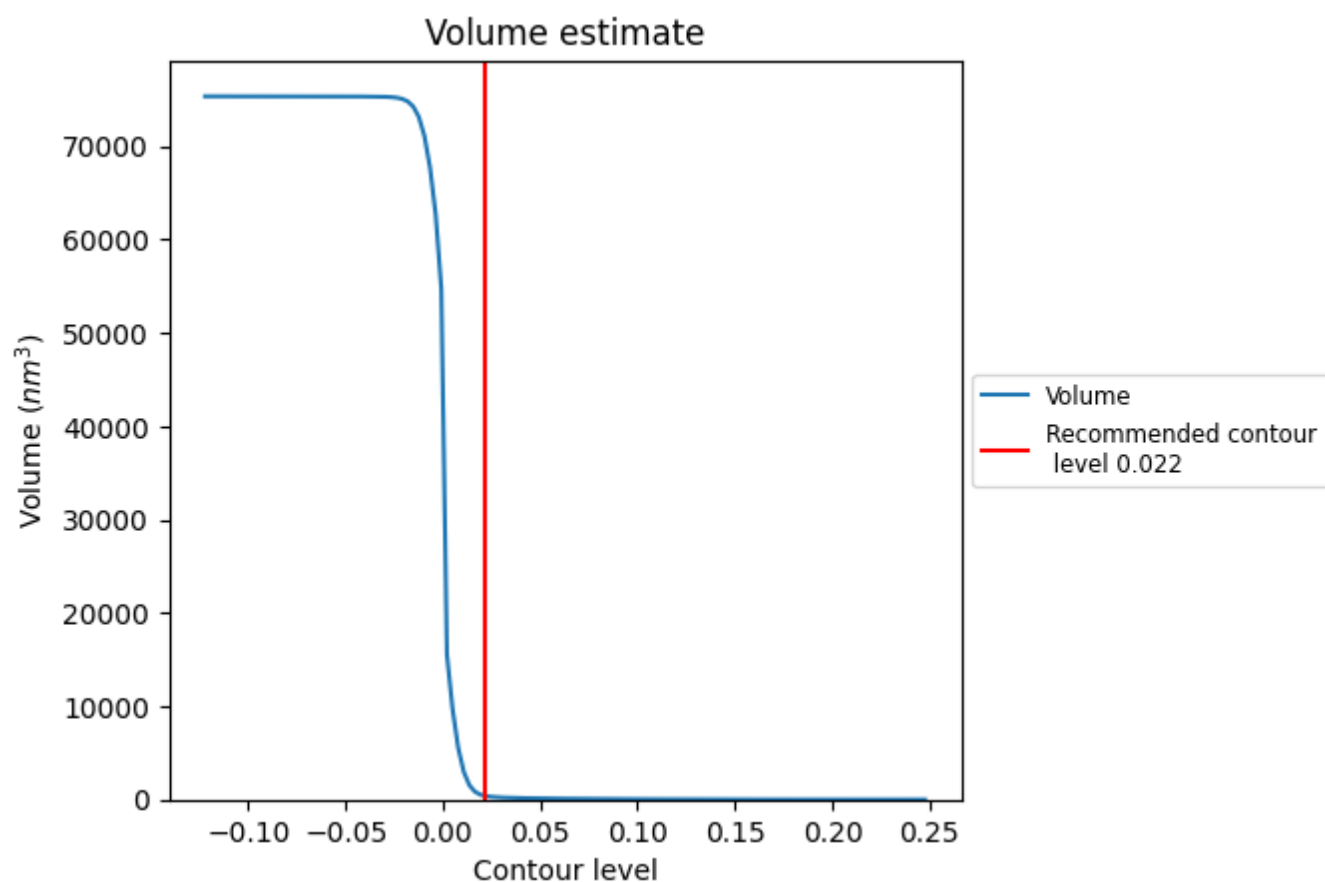
This section contains the results of statistical analysis of the map.

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.

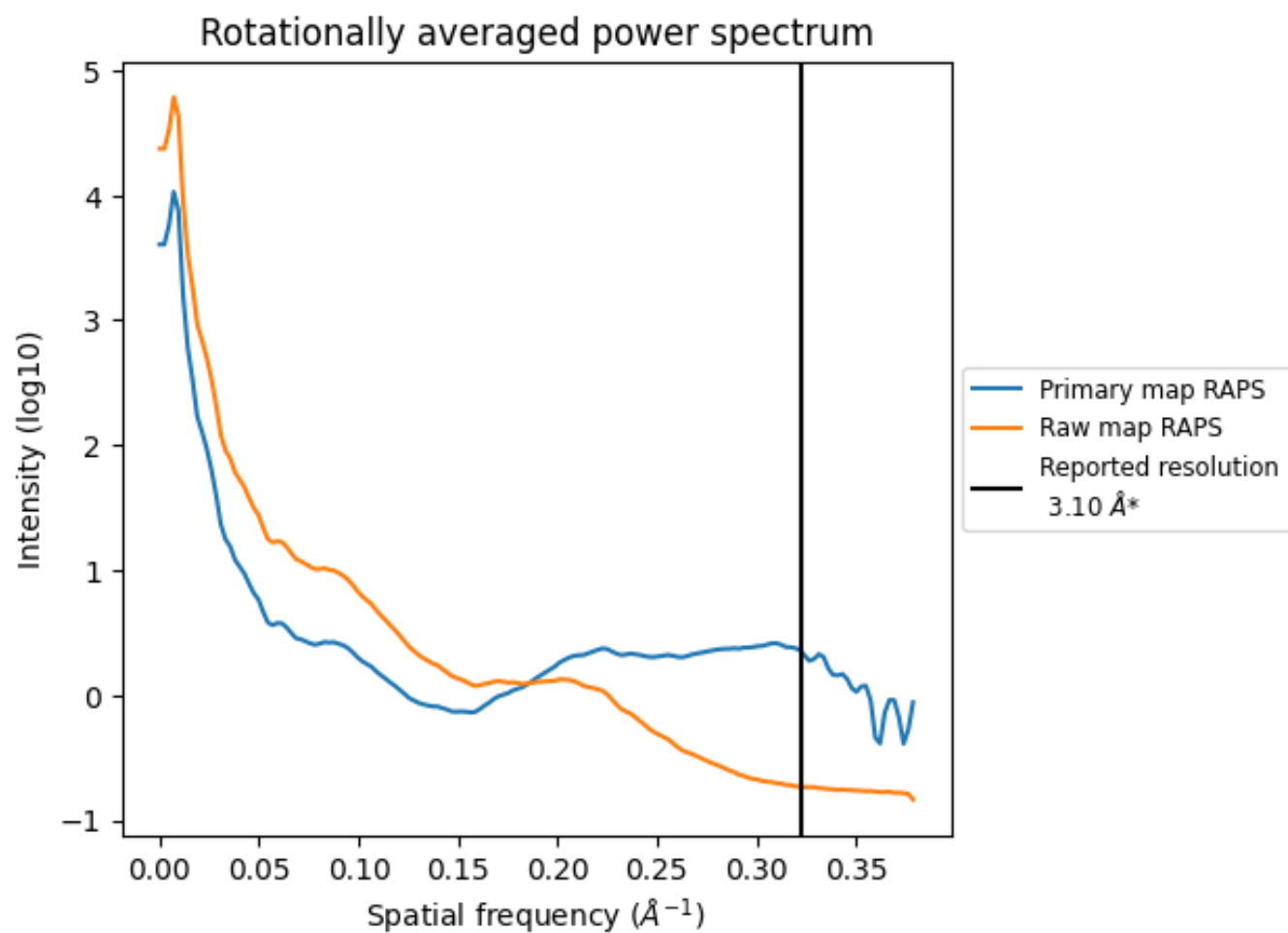
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 380 nm³; this corresponds to an approximate mass of 344 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 ⓘ

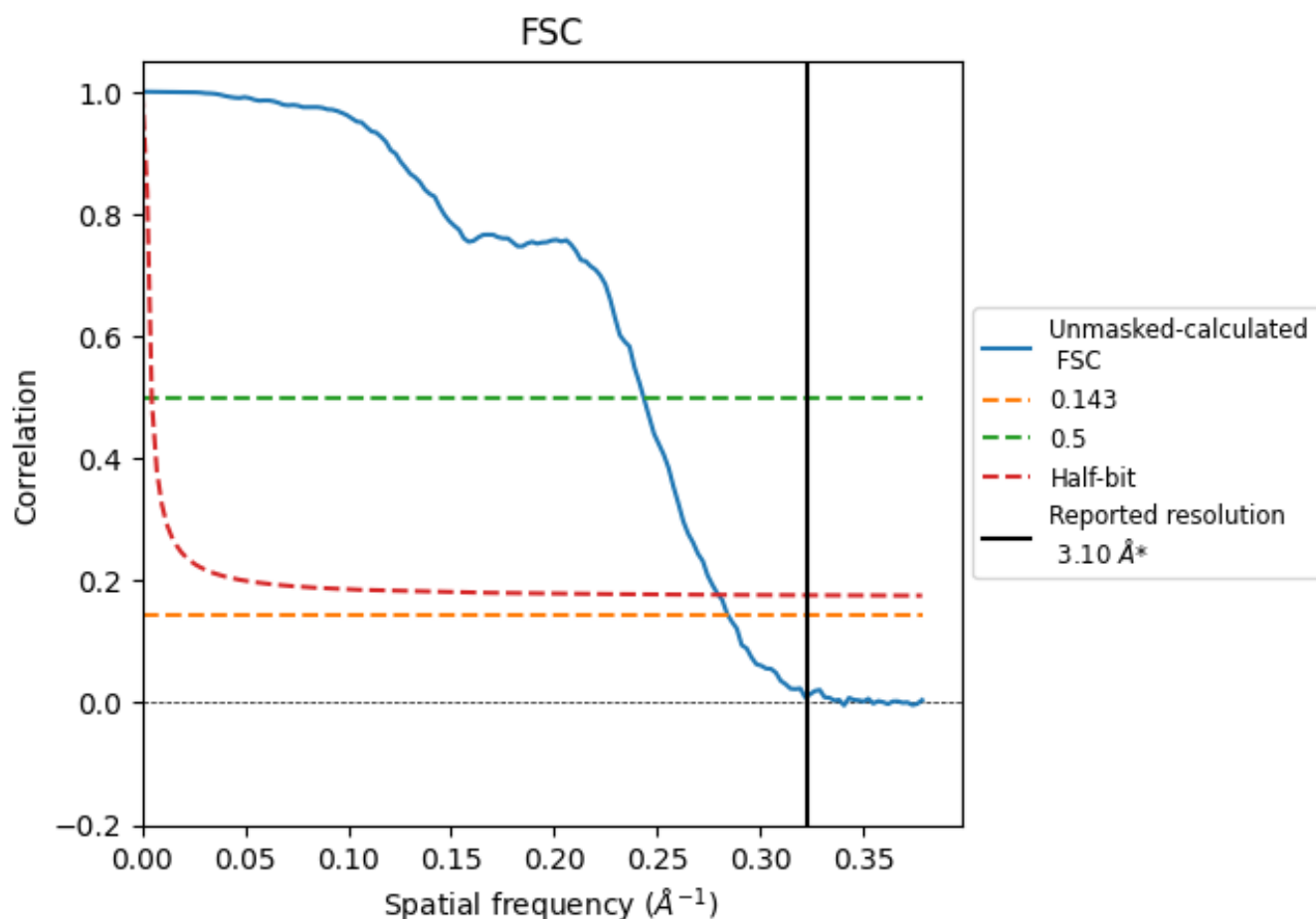


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

8.2 Resolution estimates [i](#)

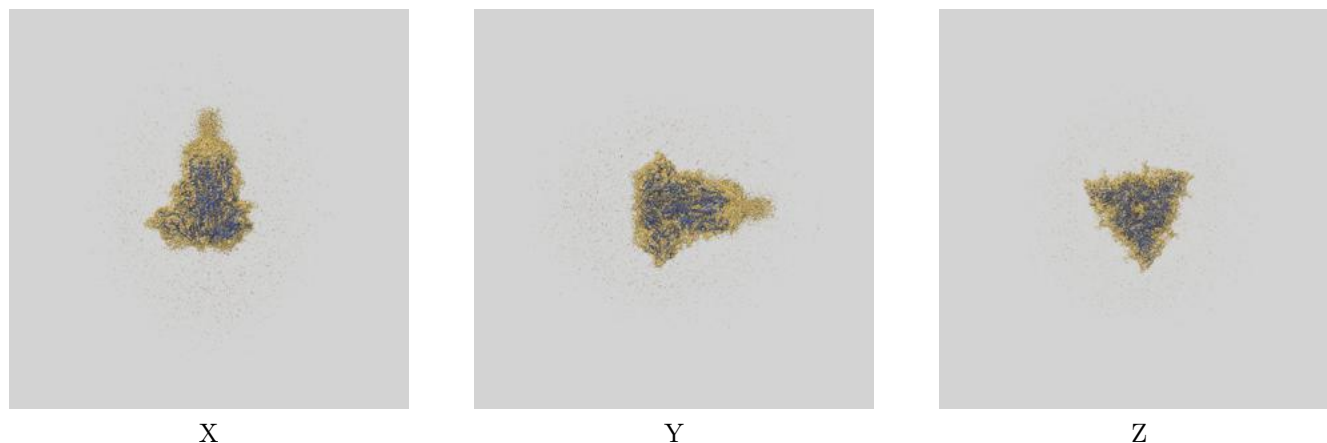
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.51	4.11	3.57

*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 3.51 differs from the reported value 3.1 by more than 10 %

9 Map-model fit [i](#)

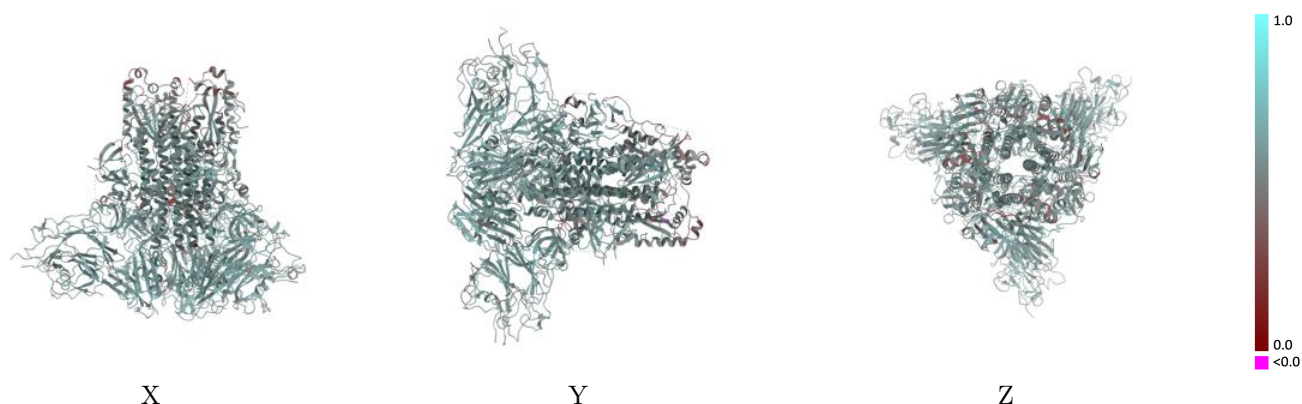
This section contains information regarding the fit between EMDB map EMD-61609 and PDB model 9JMP. Per-residue inclusion information can be found in section 3 on page 12.

9.1 Map-model overlay [i](#)



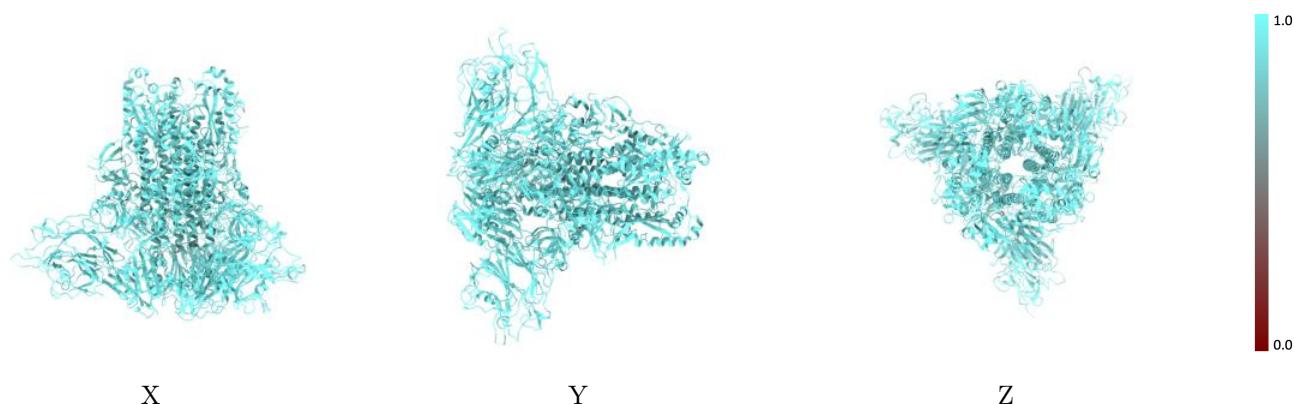
The images above show the 3D surface view of the map at the recommended contour level 0.022 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



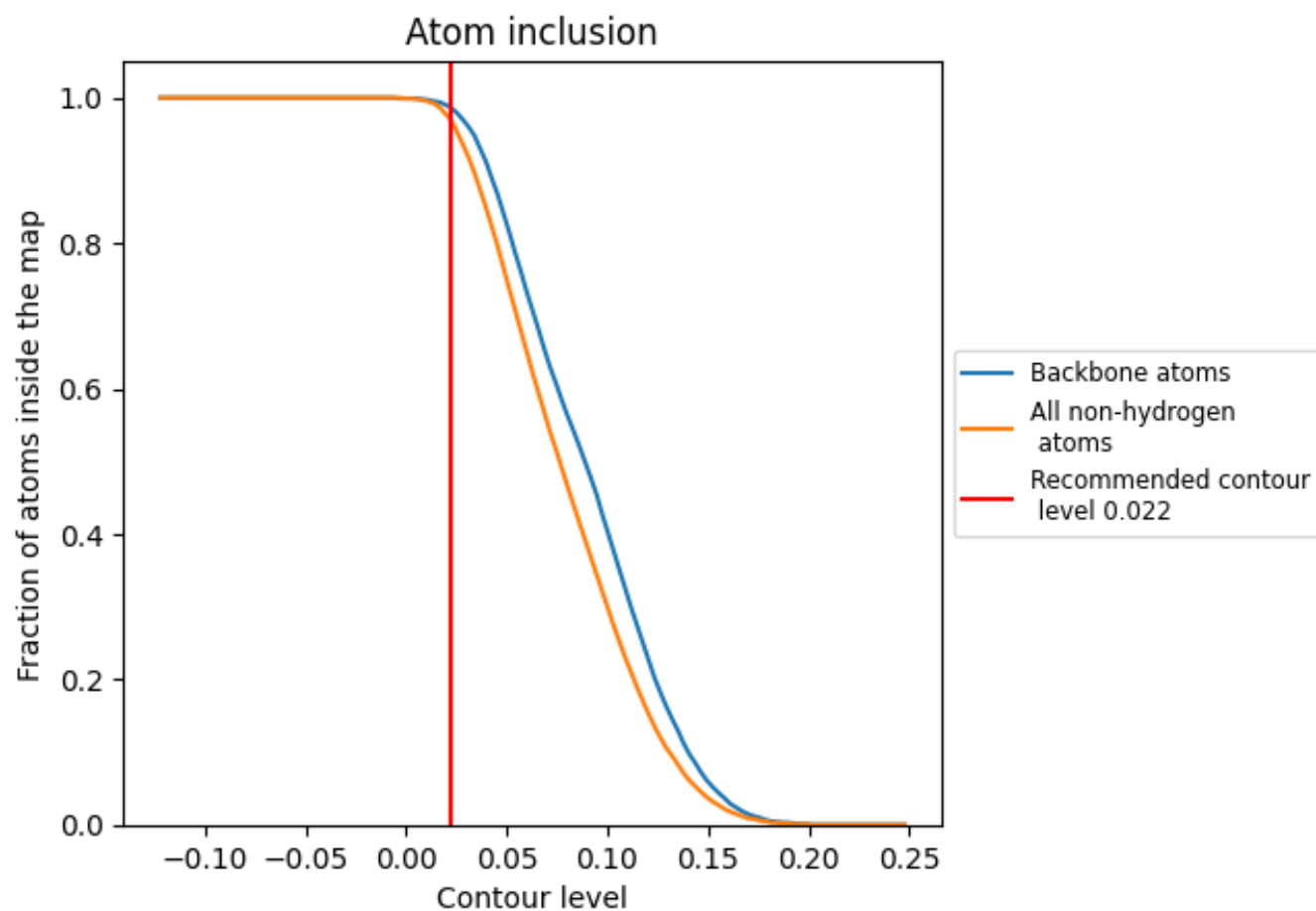
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.022).

9.4 Atom inclusion [i](#)



At the recommended contour level, 99% of all backbone atoms, 97% 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 (0.022) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.9700	<div></div> 0.5670
A	<div></div> 0.9710	<div></div> 0.5690
B	<div></div> 0.9710	<div></div> 0.5680
C	<div></div> 0.9710	<div></div> 0.5670
D	<div></div> 0.8570	<div></div> 0.5140
E	<div></div> 0.9490	<div></div> 0.5470
F	<div></div> 1.0000	<div></div> 0.5270
G	<div></div> 0.9290	<div></div> 0.4870
H	<div></div> 0.9740	<div></div> 0.5550
I	<div></div> 0.8930	<div></div> 0.4550
J	<div></div> 0.7860	<div></div> 0.3110
K	<div></div> 0.8930	<div></div> 0.5350
L	<div></div> 0.8930	<div></div> 0.5420
M	<div></div> 0.9740	<div></div> 0.5420
N	<div></div> 0.9290	<div></div> 0.4980
O	<div></div> 1.0000	<div></div> 0.5890

1.0

0.0

<0.0