



Full wwPDB EM Validation Report ⓘ

Mar 4, 2026 – 11:24 PM UTC

PDB ID : 7MXP / pdb_00007mxp
EMDB ID : EMD-24075
Title : Cryo-EM structure of NTD-directed neutralizing antibody LP5 Fab in complex with SARS-CoV-2 S2P spike
Authors : Reddem, E.R.; Casner, R.G.; Shapiro, L.
Deposited on : 2021-05-19
Resolution : 4.46 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
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

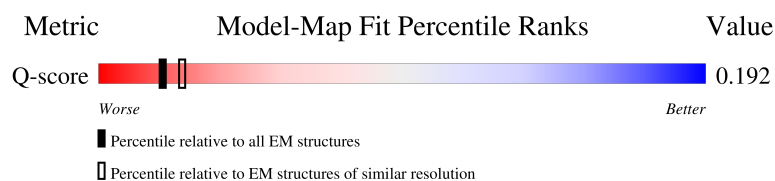
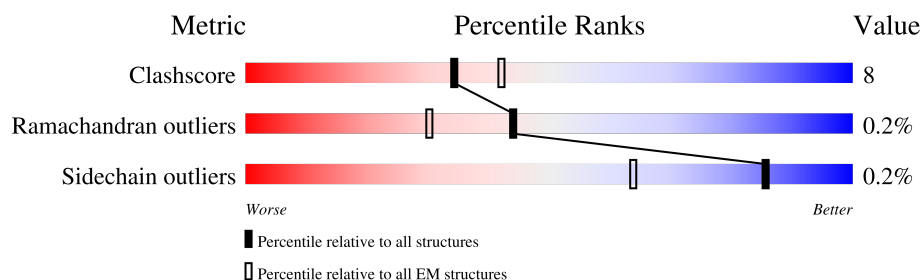
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.46 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.



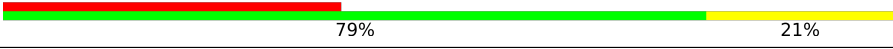
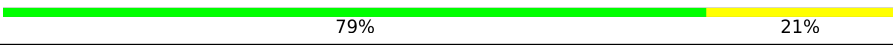
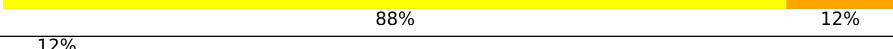
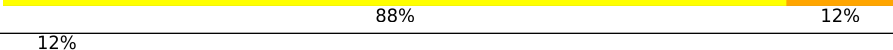
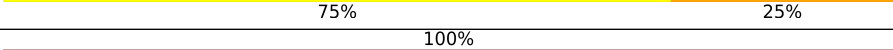
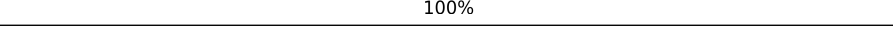
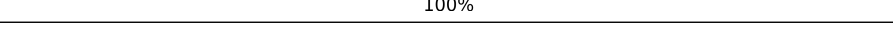
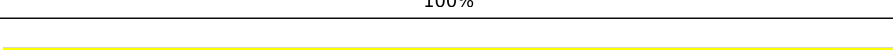
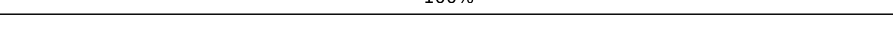
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	3001 (3.96 - 4.96)

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	1288	
1	B	1288	
1	C	1288	
2	G	119	

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Mol	Chain	Length	Quality of chain
2	H	119	
2	I	119	
3	J	107	
3	K	107	
3	L	107	
4	E	8	
4	F	8	
4	N	8	
5	D	2	
5	M	2	
5	O	2	
6	Y	3	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 31664 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	1098	Total	C	N	O	S	0	0
			8555	5458	1432	1626	39		
1	B	1101	Total	C	N	O	S	0	0
			8577	5470	1434	1634	39		
1	C	1090	Total	C	N	O	S	0	0
			8494	5416	1423	1616	39		

There are 252 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	682	GLY	ARG	conflict	UNP P0DTC2
A	683	SER	ARG	conflict	UNP P0DTC2
A	685	SER	ARG	conflict	UNP P0DTC2
A	701	VAL	ALA	variant	UNP P0DTC2
A	1209	GLY	-	expression tag	UNP P0DTC2
A	1210	SER	-	expression tag	UNP P0DTC2
A	1211	GLY	-	expression tag	UNP P0DTC2
A	1212	TYR	-	expression tag	UNP P0DTC2
A	1213	ILE	-	expression tag	UNP P0DTC2
A	1214	PRO	-	expression tag	UNP P0DTC2
A	1215	GLU	-	expression tag	UNP P0DTC2
A	1216	ALA	-	expression tag	UNP P0DTC2
A	1217	PRO	-	expression tag	UNP P0DTC2
A	1218	ARG	-	expression tag	UNP P0DTC2
A	1219	ASP	-	expression tag	UNP P0DTC2
A	1220	GLY	-	expression tag	UNP P0DTC2
A	1221	GLN	-	expression tag	UNP P0DTC2
A	1222	ALA	-	expression tag	UNP P0DTC2
A	1223	TYR	-	expression tag	UNP P0DTC2
A	1224	VAL	-	expression tag	UNP P0DTC2
A	1225	ARG	-	expression tag	UNP P0DTC2
A	1226	LYS	-	expression tag	UNP P0DTC2
A	1227	ASP	-	expression tag	UNP P0DTC2
A	1228	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1229	GLU	-	expression tag	UNP P0DTC2
A	1230	TRP	-	expression tag	UNP P0DTC2
A	1231	VAL	-	expression tag	UNP P0DTC2
A	1232	LEU	-	expression tag	UNP P0DTC2
A	1233	LEU	-	expression tag	UNP P0DTC2
A	1234	SER	-	expression tag	UNP P0DTC2
A	1235	THR	-	expression tag	UNP P0DTC2
A	1236	PHE	-	expression tag	UNP P0DTC2
A	1237	LEU	-	expression tag	UNP P0DTC2
A	1238	GLY	-	expression tag	UNP P0DTC2
A	1239	ARG	-	expression tag	UNP P0DTC2
A	1240	SER	-	expression tag	UNP P0DTC2
A	1241	LEU	-	expression tag	UNP P0DTC2
A	1242	GLU	-	expression tag	UNP P0DTC2
A	1243	VAL	-	expression tag	UNP P0DTC2
A	1244	LEU	-	expression tag	UNP P0DTC2
A	1245	PHE	-	expression tag	UNP P0DTC2
A	1246	GLN	-	expression tag	UNP P0DTC2
A	1247	GLY	-	expression tag	UNP P0DTC2
A	1248	PRO	-	expression tag	UNP P0DTC2
A	1249	GLY	-	expression tag	UNP P0DTC2
A	1250	HIS	-	expression tag	UNP P0DTC2
A	1251	HIS	-	expression tag	UNP P0DTC2
A	1252	HIS	-	expression tag	UNP P0DTC2
A	1253	HIS	-	expression tag	UNP P0DTC2
A	1254	HIS	-	expression tag	UNP P0DTC2
A	1255	HIS	-	expression tag	UNP P0DTC2
A	1256	HIS	-	expression tag	UNP P0DTC2
A	1257	HIS	-	expression tag	UNP P0DTC2
A	1258	SER	-	expression tag	UNP P0DTC2
A	1259	ALA	-	expression tag	UNP P0DTC2
A	1260	TRP	-	expression tag	UNP P0DTC2
A	1261	SER	-	expression tag	UNP P0DTC2
A	1262	HIS	-	expression tag	UNP P0DTC2
A	1263	PRO	-	expression tag	UNP P0DTC2
A	1264	GLN	-	expression tag	UNP P0DTC2
A	1265	PHE	-	expression tag	UNP P0DTC2
A	1266	GLU	-	expression tag	UNP P0DTC2
A	1267	LYS	-	expression tag	UNP P0DTC2
A	1268	GLY	-	expression tag	UNP P0DTC2
A	1269	GLY	-	expression tag	UNP P0DTC2
A	1270	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1271	SER	-	expression tag	UNP P0DTC2
A	1272	GLY	-	expression tag	UNP P0DTC2
A	1273	GLY	-	expression tag	UNP P0DTC2
A	1274	GLY	-	expression tag	UNP P0DTC2
A	1275	GLY	-	expression tag	UNP P0DTC2
A	1276	SER	-	expression tag	UNP P0DTC2
A	1277	GLY	-	expression tag	UNP P0DTC2
A	1278	GLY	-	expression tag	UNP P0DTC2
A	1279	SER	-	expression tag	UNP P0DTC2
A	1280	ALA	-	expression tag	UNP P0DTC2
A	1281	TRP	-	expression tag	UNP P0DTC2
A	1282	SER	-	expression tag	UNP P0DTC2
A	1283	HIS	-	expression tag	UNP P0DTC2
A	1284	PRO	-	expression tag	UNP P0DTC2
A	1285	GLN	-	expression tag	UNP P0DTC2
A	1286	PHE	-	expression tag	UNP P0DTC2
A	1287	GLU	-	expression tag	UNP P0DTC2
A	1288	LYS	-	expression tag	UNP P0DTC2
B	682	GLY	ARG	conflict	UNP P0DTC2
B	683	SER	ARG	conflict	UNP P0DTC2
B	685	SER	ARG	conflict	UNP P0DTC2
B	701	VAL	ALA	variant	UNP P0DTC2
B	1209	GLY	-	expression tag	UNP P0DTC2
B	1210	SER	-	expression tag	UNP P0DTC2
B	1211	GLY	-	expression tag	UNP P0DTC2
B	1212	TYR	-	expression tag	UNP P0DTC2
B	1213	ILE	-	expression tag	UNP P0DTC2
B	1214	PRO	-	expression tag	UNP P0DTC2
B	1215	GLU	-	expression tag	UNP P0DTC2
B	1216	ALA	-	expression tag	UNP P0DTC2
B	1217	PRO	-	expression tag	UNP P0DTC2
B	1218	ARG	-	expression tag	UNP P0DTC2
B	1219	ASP	-	expression tag	UNP P0DTC2
B	1220	GLY	-	expression tag	UNP P0DTC2
B	1221	GLN	-	expression tag	UNP P0DTC2
B	1222	ALA	-	expression tag	UNP P0DTC2
B	1223	TYR	-	expression tag	UNP P0DTC2
B	1224	VAL	-	expression tag	UNP P0DTC2
B	1225	ARG	-	expression tag	UNP P0DTC2
B	1226	LYS	-	expression tag	UNP P0DTC2
B	1227	ASP	-	expression tag	UNP P0DTC2
B	1228	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1229	GLU	-	expression tag	UNP P0DTC2
B	1230	TRP	-	expression tag	UNP P0DTC2
B	1231	VAL	-	expression tag	UNP P0DTC2
B	1232	LEU	-	expression tag	UNP P0DTC2
B	1233	LEU	-	expression tag	UNP P0DTC2
B	1234	SER	-	expression tag	UNP P0DTC2
B	1235	THR	-	expression tag	UNP P0DTC2
B	1236	PHE	-	expression tag	UNP P0DTC2
B	1237	LEU	-	expression tag	UNP P0DTC2
B	1238	GLY	-	expression tag	UNP P0DTC2
B	1239	ARG	-	expression tag	UNP P0DTC2
B	1240	SER	-	expression tag	UNP P0DTC2
B	1241	LEU	-	expression tag	UNP P0DTC2
B	1242	GLU	-	expression tag	UNP P0DTC2
B	1243	VAL	-	expression tag	UNP P0DTC2
B	1244	LEU	-	expression tag	UNP P0DTC2
B	1245	PHE	-	expression tag	UNP P0DTC2
B	1246	GLN	-	expression tag	UNP P0DTC2
B	1247	GLY	-	expression tag	UNP P0DTC2
B	1248	PRO	-	expression tag	UNP P0DTC2
B	1249	GLY	-	expression tag	UNP P0DTC2
B	1250	HIS	-	expression tag	UNP P0DTC2
B	1251	HIS	-	expression tag	UNP P0DTC2
B	1252	HIS	-	expression tag	UNP P0DTC2
B	1253	HIS	-	expression tag	UNP P0DTC2
B	1254	HIS	-	expression tag	UNP P0DTC2
B	1255	HIS	-	expression tag	UNP P0DTC2
B	1256	HIS	-	expression tag	UNP P0DTC2
B	1257	HIS	-	expression tag	UNP P0DTC2
B	1258	SER	-	expression tag	UNP P0DTC2
B	1259	ALA	-	expression tag	UNP P0DTC2
B	1260	TRP	-	expression tag	UNP P0DTC2
B	1261	SER	-	expression tag	UNP P0DTC2
B	1262	HIS	-	expression tag	UNP P0DTC2
B	1263	PRO	-	expression tag	UNP P0DTC2
B	1264	GLN	-	expression tag	UNP P0DTC2
B	1265	PHE	-	expression tag	UNP P0DTC2
B	1266	GLU	-	expression tag	UNP P0DTC2
B	1267	LYS	-	expression tag	UNP P0DTC2
B	1268	GLY	-	expression tag	UNP P0DTC2
B	1269	GLY	-	expression tag	UNP P0DTC2
B	1270	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1271	SER	-	expression tag	UNP P0DTC2
B	1272	GLY	-	expression tag	UNP P0DTC2
B	1273	GLY	-	expression tag	UNP P0DTC2
B	1274	GLY	-	expression tag	UNP P0DTC2
B	1275	GLY	-	expression tag	UNP P0DTC2
B	1276	SER	-	expression tag	UNP P0DTC2
B	1277	GLY	-	expression tag	UNP P0DTC2
B	1278	GLY	-	expression tag	UNP P0DTC2
B	1279	SER	-	expression tag	UNP P0DTC2
B	1280	ALA	-	expression tag	UNP P0DTC2
B	1281	TRP	-	expression tag	UNP P0DTC2
B	1282	SER	-	expression tag	UNP P0DTC2
B	1283	HIS	-	expression tag	UNP P0DTC2
B	1284	PRO	-	expression tag	UNP P0DTC2
B	1285	GLN	-	expression tag	UNP P0DTC2
B	1286	PHE	-	expression tag	UNP P0DTC2
B	1287	GLU	-	expression tag	UNP P0DTC2
B	1288	LYS	-	expression tag	UNP P0DTC2
C	682	GLY	ARG	conflict	UNP P0DTC2
C	683	SER	ARG	conflict	UNP P0DTC2
C	685	SER	ARG	conflict	UNP P0DTC2
C	701	VAL	ALA	variant	UNP P0DTC2
C	1209	GLY	-	expression tag	UNP P0DTC2
C	1210	SER	-	expression tag	UNP P0DTC2
C	1211	GLY	-	expression tag	UNP P0DTC2
C	1212	TYR	-	expression tag	UNP P0DTC2
C	1213	ILE	-	expression tag	UNP P0DTC2
C	1214	PRO	-	expression tag	UNP P0DTC2
C	1215	GLU	-	expression tag	UNP P0DTC2
C	1216	ALA	-	expression tag	UNP P0DTC2
C	1217	PRO	-	expression tag	UNP P0DTC2
C	1218	ARG	-	expression tag	UNP P0DTC2
C	1219	ASP	-	expression tag	UNP P0DTC2
C	1220	GLY	-	expression tag	UNP P0DTC2
C	1221	GLN	-	expression tag	UNP P0DTC2
C	1222	ALA	-	expression tag	UNP P0DTC2
C	1223	TYR	-	expression tag	UNP P0DTC2
C	1224	VAL	-	expression tag	UNP P0DTC2
C	1225	ARG	-	expression tag	UNP P0DTC2
C	1226	LYS	-	expression tag	UNP P0DTC2
C	1227	ASP	-	expression tag	UNP P0DTC2
C	1228	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1229	GLU	-	expression tag	UNP P0DTC2
C	1230	TRP	-	expression tag	UNP P0DTC2
C	1231	VAL	-	expression tag	UNP P0DTC2
C	1232	LEU	-	expression tag	UNP P0DTC2
C	1233	LEU	-	expression tag	UNP P0DTC2
C	1234	SER	-	expression tag	UNP P0DTC2
C	1235	THR	-	expression tag	UNP P0DTC2
C	1236	PHE	-	expression tag	UNP P0DTC2
C	1237	LEU	-	expression tag	UNP P0DTC2
C	1238	GLY	-	expression tag	UNP P0DTC2
C	1239	ARG	-	expression tag	UNP P0DTC2
C	1240	SER	-	expression tag	UNP P0DTC2
C	1241	LEU	-	expression tag	UNP P0DTC2
C	1242	GLU	-	expression tag	UNP P0DTC2
C	1243	VAL	-	expression tag	UNP P0DTC2
C	1244	LEU	-	expression tag	UNP P0DTC2
C	1245	PHE	-	expression tag	UNP P0DTC2
C	1246	GLN	-	expression tag	UNP P0DTC2
C	1247	GLY	-	expression tag	UNP P0DTC2
C	1248	PRO	-	expression tag	UNP P0DTC2
C	1249	GLY	-	expression tag	UNP P0DTC2
C	1250	HIS	-	expression tag	UNP P0DTC2
C	1251	HIS	-	expression tag	UNP P0DTC2
C	1252	HIS	-	expression tag	UNP P0DTC2
C	1253	HIS	-	expression tag	UNP P0DTC2
C	1254	HIS	-	expression tag	UNP P0DTC2
C	1255	HIS	-	expression tag	UNP P0DTC2
C	1256	HIS	-	expression tag	UNP P0DTC2
C	1257	HIS	-	expression tag	UNP P0DTC2
C	1258	SER	-	expression tag	UNP P0DTC2
C	1259	ALA	-	expression tag	UNP P0DTC2
C	1260	TRP	-	expression tag	UNP P0DTC2
C	1261	SER	-	expression tag	UNP P0DTC2
C	1262	HIS	-	expression tag	UNP P0DTC2
C	1263	PRO	-	expression tag	UNP P0DTC2
C	1264	GLN	-	expression tag	UNP P0DTC2
C	1265	PHE	-	expression tag	UNP P0DTC2
C	1266	GLU	-	expression tag	UNP P0DTC2
C	1267	LYS	-	expression tag	UNP P0DTC2
C	1268	GLY	-	expression tag	UNP P0DTC2
C	1269	GLY	-	expression tag	UNP P0DTC2
C	1270	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1271	SER	-	expression tag	UNP P0DTC2
C	1272	GLY	-	expression tag	UNP P0DTC2
C	1273	GLY	-	expression tag	UNP P0DTC2
C	1274	GLY	-	expression tag	UNP P0DTC2
C	1275	GLY	-	expression tag	UNP P0DTC2
C	1276	SER	-	expression tag	UNP P0DTC2
C	1277	GLY	-	expression tag	UNP P0DTC2
C	1278	GLY	-	expression tag	UNP P0DTC2
C	1279	SER	-	expression tag	UNP P0DTC2
C	1280	ALA	-	expression tag	UNP P0DTC2
C	1281	TRP	-	expression tag	UNP P0DTC2
C	1282	SER	-	expression tag	UNP P0DTC2
C	1283	HIS	-	expression tag	UNP P0DTC2
C	1284	PRO	-	expression tag	UNP P0DTC2
C	1285	GLN	-	expression tag	UNP P0DTC2
C	1286	PHE	-	expression tag	UNP P0DTC2
C	1287	GLU	-	expression tag	UNP P0DTC2
C	1288	LYS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called LP5 Fab Heavy Chain.

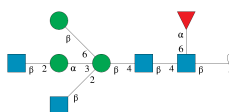
Mol	Chain	Residues	Atoms					AltConf	Trace
2	G	119	Total	C	N	O	S	0	0
			936	597	159	176	4		
2	I	119	Total	C	N	O	S	0	0
			936	597	159	176	4		
2	H	119	Total	C	N	O	S	0	0
			936	597	159	176	4		

- Molecule 3 is a protein called LP5 Fab Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	J	107	Total	C	N	O	S	0	0
			806	506	131	167	2		
3	K	107	Total	C	N	O	S	0	0
			806	506	131	167	2		
3	L	107	Total	C	N	O	S	0	0
			806	506	131	167	2		

- Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)][beta-D-mannopyranose-(1-6)][beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1

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



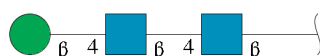
Mol	Chain	Residues	Atoms				AltConf	Trace
4	E	8	Total	C	N	O	0	0
			99	56	4	39		
4	F	8	Total	C	N	O	0	0
			99	56	4	39		
4	N	8	Total	C	N	O	0	0
			99	56	4	39		

- Molecule 5 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
5	D	2	Total	C	N	O	0	0
			28	16	2	10		
5	M	2	Total	C	N	O	0	0
			28	16	2	10		
5	O	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 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
6	Y	3	Total	C	N	O	0	0
			39	22	2	15		

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



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

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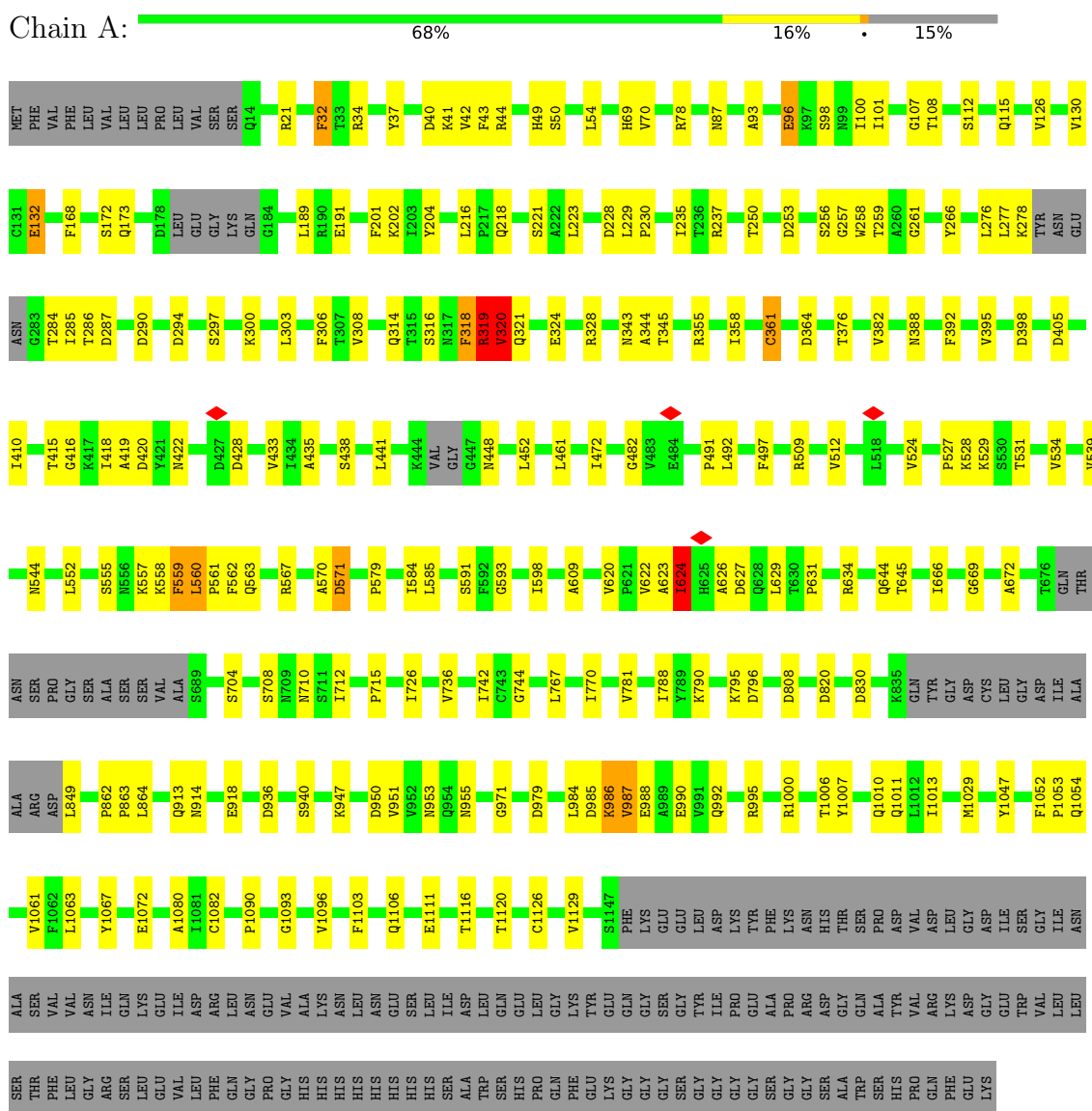
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Mol	Chain	Residues	Atoms				AltConf
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	B	1	Total	C	N	O	0
			14	8	1	5	
7	C	1	Total	C	N	O	0
			14	8	1	5	
7	C	1	Total	C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

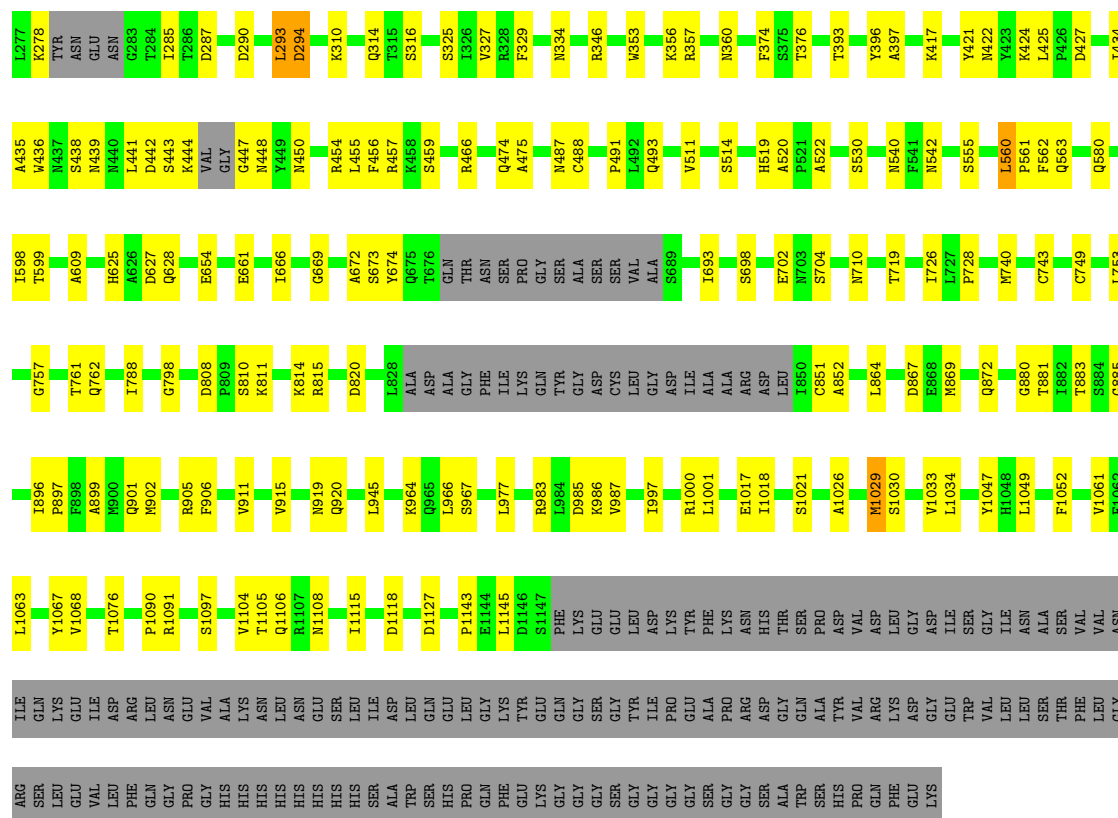
3 Residue-property plots

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.

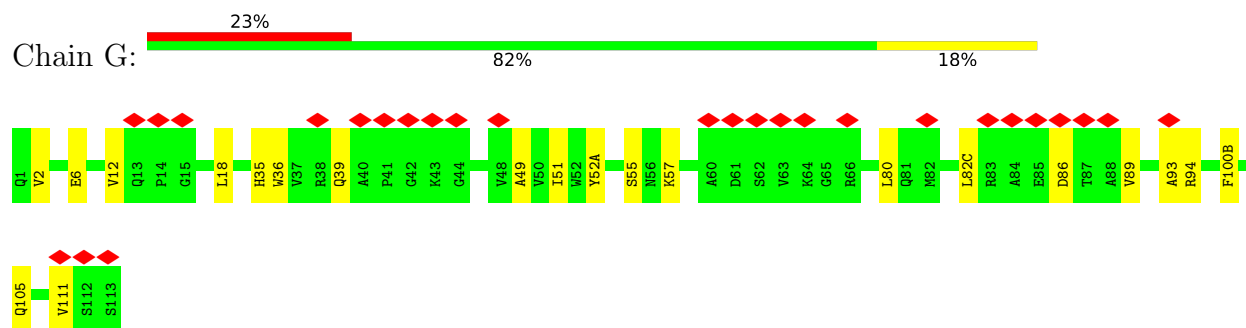
• Molecule 1: Spike glycoprotein



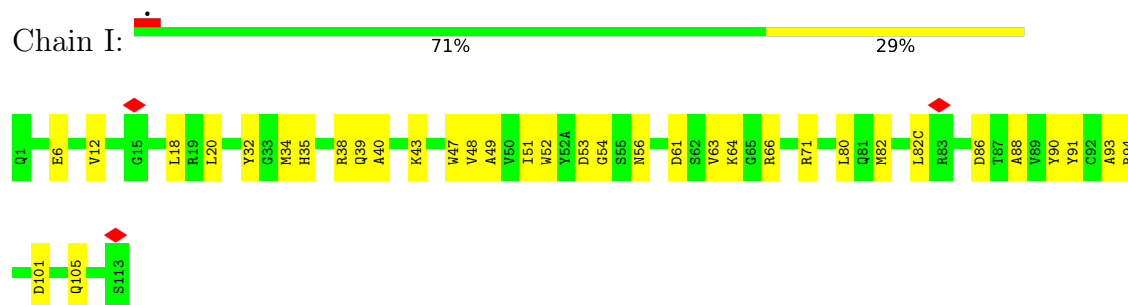




- Molecule 2: LP5 Fab Heavy Chain

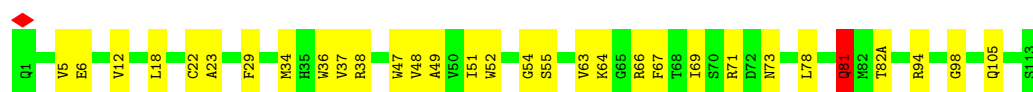


- Molecule 2: LP5 Fab Heavy Chain

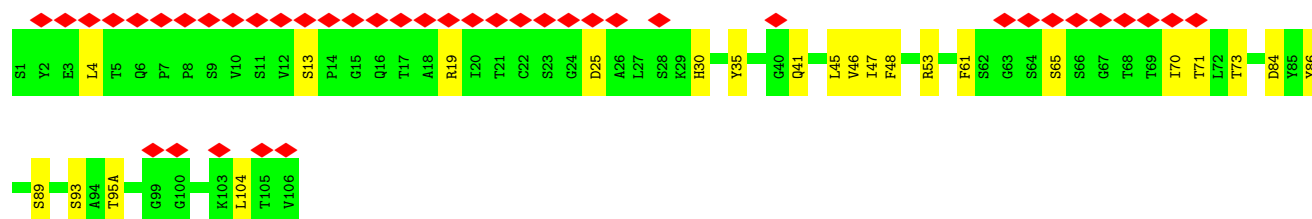
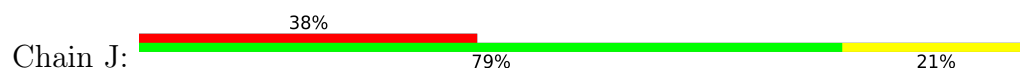


- Molecule 2: LP5 Fab Heavy Chain

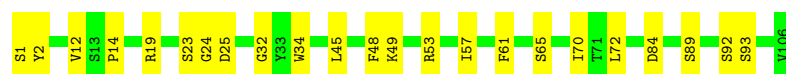
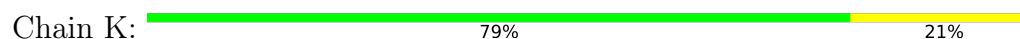




- Molecule 3: LP5 Fab Light Chain



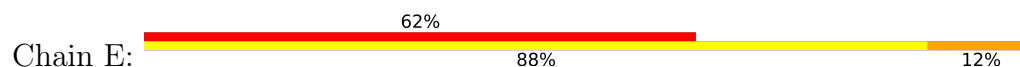
- Molecule 3: LP5 Fab Light Chain



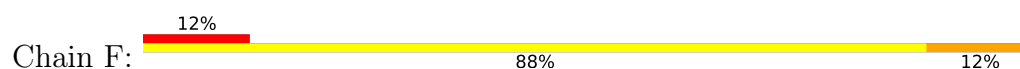
- Molecule 3: LP5 Fab Light Chain



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)][beta-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)][beta-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

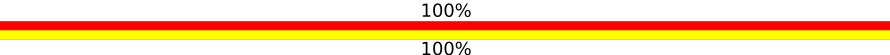


- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)][beta-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N: 

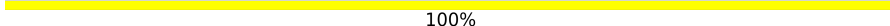


- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D: 

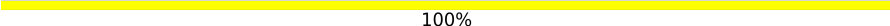


- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M: 

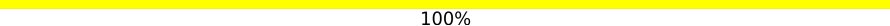


- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O: 



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

Chain Y: 



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	96531	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 ($e^-/\text{\AA}^2$)	42	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.801	Depositor
Minimum map value	-0.161	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.028	Depositor
Recommended contour level	0.15	Depositor
Map size (Å)	479.36002, 479.36002, 479.36002	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.07, 1.07, 1.07	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, FUC, BMA, MAN

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	0.41	0/8753	0.87	29/11913 (0.2%)
1	B	0.46	2/8776 (0.0%)	0.88	15/11945 (0.1%)
1	C	0.41	1/8691 (0.0%)	0.85	15/11830 (0.1%)
2	G	0.44	0/961	0.93	0/1302
2	H	0.41	0/961	0.88	1/1302 (0.1%)
2	I	0.43	0/961	0.89	1/1302 (0.1%)
3	J	0.41	0/827	0.82	0/1128
3	K	0.37	0/827	0.86	2/1128 (0.2%)
3	L	0.38	0/827	0.87	0/1128
All	All	0.42	3/31584 (0.0%)	0.87	63/42978 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	8
1	B	0	9
1	C	0	10
2	G	0	1
3	J	0	2
3	L	0	1
All	All	0	31

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	986	LYS	C-N	10.90	1.45	1.34
1	C	986	LYS	C-N	9.15	1.45	1.34
1	B	289	VAL	N-CA	5.19	1.52	1.46

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	559	PHE	CA-CB-CG	7.34	121.14	113.80
1	B	42	VAL	CA-CB-CG1	7.16	122.57	110.40
1	A	624	ILE	CG1-CB-CG2	-7.03	89.60	110.70
1	C	986	LYS	N-CA-C	6.70	119.41	111.71
1	B	744	GLY	CA-C-N	6.47	131.59	122.08
1	B	744	GLY	C-N-CA	6.47	131.59	122.08
1	A	132	GLU	CA-CB-CG	6.42	126.93	114.10
1	C	1017	GLU	N-CA-CB	6.41	120.22	110.28
1	B	359	SER	CA-C-N	6.40	131.48	122.46
1	B	359	SER	C-N-CA	6.40	131.48	122.46
1	A	570	ALA	CA-C-N	6.33	137.37	125.66
1	A	570	ALA	C-N-CA	6.33	137.37	125.66
2	H	81	GLN	CA-CB-CG	6.32	126.75	114.10
1	B	631	PRO	CA-N-CD	-6.21	103.31	112.00
1	C	987	VAL	N-CA-C	6.13	118.71	111.05
1	A	392	PHE	CA-C-N	6.13	133.24	121.54
1	A	392	PHE	C-N-CA	6.13	133.24	121.54
1	B	290	ASP	CA-C-N	6.11	133.03	122.64
1	B	290	ASP	C-N-CA	6.11	133.03	122.64
1	B	42	VAL	N-CA-CB	-6.08	101.20	111.23
1	A	629	LEU	CA-C-N	6.01	128.75	120.39
1	A	629	LEU	C-N-CA	6.01	128.75	120.39
1	B	986	LYS	N-CA-C	5.98	117.80	111.28
1	A	559	PHE	CA-C-N	5.98	136.39	121.80
1	A	559	PHE	C-N-CA	5.98	136.39	121.80
1	B	1005	GLN	N-CA-CB	5.90	119.43	110.28
2	I	43	LYS	CA-CB-CG	5.89	125.87	114.10
1	A	558	LYS	CA-CB-CG	5.88	125.85	114.10
1	C	519	HIS	CA-C-N	5.86	129.01	120.51
1	C	519	HIS	C-N-CA	5.86	129.01	120.51
1	B	289	VAL	N-CA-C	5.78	116.49	108.23
1	C	276	LEU	CA-CB-CG	5.61	135.92	116.30
1	C	1143	PRO	CA-N-CD	-5.57	104.20	112.00
1	A	343	ASN	N-CA-C	5.51	117.28	111.28
1	C	1017	GLU	CA-CB-CG	5.48	125.07	114.10
3	K	49	LYS	CA-C-N	5.47	131.99	121.54
3	K	49	LYS	C-N-CA	5.47	131.99	121.54
1	A	744	GLY	CA-C-N	5.42	129.83	122.07
1	A	744	GLY	C-N-CA	5.42	129.83	122.07
1	C	555	SER	CA-C-N	5.42	131.90	121.54
1	C	555	SER	C-N-CA	5.42	131.90	121.54
1	B	570	ALA	CA-C-N	5.42	131.89	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	570	ALA	C-N-CA	5.42	131.89	121.54
1	A	410	ILE	N-CA-C	-5.41	108.00	113.47
1	A	410	ILE	CB-CA-C	5.38	116.43	110.62
1	C	258	TRP	CA-C-N	5.25	130.27	122.82
1	C	258	TRP	C-N-CA	5.25	130.27	122.82
1	A	620	VAL	N-CA-C	5.21	120.17	112.61
1	A	258	TRP	CA-C-N	5.16	130.15	122.82
1	A	258	TRP	C-N-CA	5.16	130.15	122.82
1	A	560	LEU	N-CA-CB	-5.14	101.21	110.37
1	A	87	ASN	CA-C-N	5.11	129.59	122.08
1	A	87	ASN	C-N-CA	5.11	129.59	122.08
1	A	259	THR	CA-CB-CG2	5.09	119.16	110.50
1	A	979	ASP	CA-CB-CG	5.08	117.68	112.60
1	B	1005	GLN	CA-CB-CG	5.08	124.25	114.10
1	C	1029	MET	CB-CG-SD	5.07	127.91	112.70
1	A	979	ASP	CB-CG-OD2	5.03	129.98	118.40
1	C	293	LEU	CA-C-N	5.02	134.06	121.80
1	C	293	LEU	C-N-CA	5.02	134.06	121.80
1	A	461	LEU	CA-C-N	5.00	127.34	120.39
1	A	461	LEU	C-N-CA	5.00	127.34	120.39
1	A	571	ASP	CA-CB-CG	5.00	117.60	112.60

There are no chirality outliers.

All (31) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	21	ARG	Peptide
1	A	290	ASP	Peptide
1	A	32	PHE	Peptide
1	A	344	ALA	Peptide
1	A	361	CYS	Peptide
1	A	561	PRO	Peptide
1	A	624	ILE	Peptide
1	A	736	VAL	Peptide
1	B	15	CYS	Peptide
1	B	21	ARG	Peptide
1	B	248	TYR	Peptide
1	B	285	ILE	Mainchain
1	B	39	PRO	Peptide
1	B	590	CYS	Peptide
1	B	617	CYS	Peptide
1	B	631	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	B	855	PHE	Peptide
1	C	1145	LEU	Peptide
1	C	21	ARG	Peptide
1	C	290	ASP	Peptide
1	C	293	LEU	Peptide
1	C	294	ASP	Peptide
1	C	32	PHE	Peptide
1	C	520	ALA	Peptide
1	C	560	LEU	Peptide
1	C	562	PHE	Peptide
1	C	740	MET	Peptide
2	G	80	LEU	Peptide
3	J	13	SER	Peptide
3	J	30	HIS	Peptide
3	L	13	SER	Peptide

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	8555	0	8328	142	0
1	B	8577	0	8335	174	0
1	C	8494	0	8259	125	0
2	G	936	0	886	13	0
2	H	936	0	886	16	0
2	I	936	0	886	23	0
3	J	806	0	762	17	0
3	K	806	0	762	13	0
3	L	806	0	762	19	0
4	E	99	0	85	1	0
4	F	99	0	85	1	0
4	N	99	0	85	1	0
5	D	28	0	25	0	0
5	M	28	0	25	0	0
5	O	28	0	25	0	0
6	Y	39	0	34	0	0
7	A	126	0	117	0	0
7	B	140	0	130	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	C	126	0	117	1	0
All	All	31664	0	30594	503	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (503) 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:275:PHE:HA	1:B:289:VAL:C	1.97	0.87
1:A:319:ARG:HB3	1:A:591:SER:O	1.81	0.81
1:A:320:VAL:H	1:A:591:SER:CB	1.95	0.80
1:C:1029:MET:O	1:C:1033:VAL:HB	1.84	0.77
1:B:275:PHE:HA	1:B:289:VAL:HA	1.67	0.75
1:B:1029:MET:O	1:B:1033:VAL:HB	1.88	0.74
1:B:275:PHE:HA	1:B:289:VAL:CA	2.20	0.71
1:B:277:LEU:HA	1:B:287:ASP:HA	1.70	0.71
1:B:1028:LYS:O	1:B:1032:CYS:HB2	1.91	0.71
1:B:916:LEU:O	1:B:920:GLN:HB3	1.93	0.68
1:B:278:LYS:HB3	1:B:286:THR:H	1.58	0.67
1:B:276:LEU:O	1:B:288:ALA:N	2.28	0.67
1:B:206:LYS:HZ2	1:B:222:ALA:H	1.43	0.67
1:A:320:VAL:H	1:A:591:SER:HB3	1.58	0.66
1:A:320:VAL:H	1:A:591:SER:HB2	1.59	0.66
1:A:562:PHE:HB2	1:B:225:PRO:HG2	1.80	0.64
1:A:557:LYS:H	1:B:43:PHE:HB3	1.62	0.64
1:A:986:LYS:O	1:A:987:VAL:C	2.41	0.63
1:B:1015:ALA:O	1:B:1019:ARG:HB3	1.99	0.62
1:A:1047:TYR:HB2	1:A:1067:TYR:HB3	1.81	0.61
1:B:275:PHE:CA	1:B:289:VAL:HA	2.32	0.60
1:A:986:LYS:O	1:A:990:GLU:N	2.35	0.59
1:B:275:PHE:HB2	1:B:288:ALA:O	2.02	0.59
1:B:93:ALA:HB3	1:B:266:TYR:HB2	1.85	0.59
1:B:275:PHE:HA	1:B:290:ASP:N	2.16	0.59
1:B:276:LEU:H	1:B:289:VAL:N	2.01	0.59
2:G:35:HIS:HB2	2:G:93:ALA:HB3	1.83	0.59
2:H:81:GLN:NE2	2:H:82(A):THR:OG1	2.32	0.58
1:B:189:LEU:HD22	1:B:217:PRO:HD2	1.85	0.58
1:C:1106:GLN:HG2	1:C:1108:ASN:H	1.68	0.57
1:C:376:THR:HB	1:C:435:ALA:HB3	1.85	0.57
1:B:93:ALA:HB1	1:B:189:LEU:HD11	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:69:HIS:HB3	1:C:78:ARG:HD2	1.86	0.57
1:B:992:GLN:HA	1:B:995:ARG:HH21	1.70	0.57
1:C:710:ASN:HB3	7:C:1306:NAG:H82	1.85	0.57
3:J:46:VAL:HG23	3:J:47:ILE:HG22	1.85	0.57
1:A:70:VAL:HG21	1:A:261:GLY:H	1.70	0.57
1:C:55:PHE:HB2	1:C:273:ARG:HB3	1.88	0.56
3:L:12:VAL:HG21	3:L:18:ALA:H	1.69	0.56
1:A:388:ASN:HB2	1:A:527:PRO:HD2	1.88	0.56
1:A:448:ASN:HB3	1:A:497:PHE:HB2	1.87	0.56
2:I:35:HIS:HB2	2:I:93:ALA:HB3	1.88	0.56
1:B:36:VAL:HG11	1:B:287:ASP:HB2	1.88	0.56
1:B:878:LEU:O	1:B:882:ILE:HB	2.05	0.56
1:B:773:GLU:HA	1:B:776:LYS:HG2	1.88	0.55
1:C:851:CYS:SG	1:C:852:ALA:N	2.79	0.55
2:I:66:ARG:HD2	2:I:82(C):LEU:HA	1.89	0.55
1:A:1054:GLN:HB2	1:A:1061:VAL:HB	1.89	0.54
1:C:1047:TYR:HB2	1:C:1067:TYR:HB3	1.89	0.54
1:C:1076:THR:HB	1:C:1097:SER:HB3	1.89	0.54
2:G:6:GLU:O	2:G:105:GLN:NE2	2.40	0.54
1:A:50:SER:HA	1:A:276:LEU:HA	1.89	0.54
1:A:93:ALA:HB1	1:A:189:LEU:HD11	1.89	0.54
1:A:126:VAL:HB	1:A:172:SER:HB3	1.89	0.54
1:B:275:PHE:HB2	1:B:289:VAL:HA	1.89	0.54
1:A:324:GLU:HB2	1:A:539:VAL:HG12	1.88	0.54
1:B:214:ARG:HH12	1:B:262:ALA:HA	1.72	0.54
1:C:356:LYS:HB3	1:C:397:ALA:HB3	1.89	0.54
1:B:710:ASN:HB3	7:B:1305:NAG:H82	1.89	0.54
1:C:200:TYR:HB3	1:C:228:ASP:HB2	1.90	0.54
3:L:53:ARG:NH1	3:L:61:PHE:O	2.41	0.54
1:C:997:ILE:O	1:C:1001:LEU:HB2	2.07	0.54
1:B:92:PHE:HB2	1:B:192:PHE:HB2	1.90	0.53
3:K:53:ARG:HB3	3:K:57:ILE:HG21	1.90	0.53
1:B:186:PHE:HB2	1:B:210:ILE:HB	1.91	0.53
1:C:50:SER:HA	1:C:276:LEU:HA	1.90	0.53
1:C:75:GLY:HA2	1:C:78:ARG:HH11	1.73	0.53
1:A:49:HIS:HB3	1:A:277:LEU:HD12	1.91	0.53
2:H:64:LYS:O	2:H:66:ARG:NH1	2.41	0.53
1:B:953:ASN:O	1:B:957:GLN:HB2	2.08	0.53
2:H:6:GLU:HA	2:H:22:CYS:HA	1.90	0.53
1:B:560:LEU:O	1:B:577:ARG:NH2	2.42	0.53
1:A:328:ARG:NH2	1:A:531:THR:O	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:474:GLN:NE2	1:B:488:CYS:SG	2.82	0.53
1:B:867:ASP:OD1	1:B:867:ASP:N	2.42	0.53
1:B:314:GLN:NE2	1:B:316:SER:O	2.41	0.53
1:C:195:LYS:HE2	1:C:197:ILE:HD11	1.91	0.53
1:A:704:SER:HA	1:B:790:LYS:HB3	1.91	0.53
3:L:65:SER:HA	3:L:70:ILE:HA	1.91	0.52
1:B:314:GLN:OE1	1:B:613:GLN:NE2	2.42	0.52
1:A:1007:TYR:O	1:A:1011:GLN:NE2	2.42	0.52
1:B:274:THR:O	1:B:290:ASP:N	2.43	0.52
1:B:275:PHE:CB	1:B:289:VAL:HA	2.40	0.52
1:B:739:THR:HA	1:B:742:ILE:HG22	1.91	0.52
1:B:905:ARG:NH1	1:B:1049:LEU:O	2.43	0.52
3:J:84:ASP:OD1	3:J:84:ASP:N	2.43	0.52
1:C:1104:VAL:HG23	1:C:1115:ILE:HG12	1.91	0.52
2:I:20:LEU:HB2	2:I:80:LEU:HD12	1.90	0.52
3:L:53:ARG:NH1	3:L:73:THR:OG1	2.42	0.52
1:A:115:GLN:HA	1:A:132:GLU:HB3	1.91	0.52
1:B:776:LYS:HD2	1:B:1019:ARG:HH22	1.73	0.52
1:C:104:TRP:HA	1:C:240:THR:HA	1.91	0.52
1:C:131:CYS:SG	1:C:132:GLU:N	2.83	0.52
1:B:518:LEU:HD12	1:C:983:ARG:HH21	1.74	0.51
1:A:364:ASP:OD1	1:A:388:ASN:ND2	2.44	0.51
3:J:65:SER:H	3:J:70:ILE:HG23	1.74	0.51
1:A:172:SER:OG	1:A:173:GLN:N	2.43	0.51
1:B:951:VAL:O	1:B:955:ASN:ND2	2.43	0.51
2:I:82:MET:HE2	2:I:82(C):LEU:HD22	1.93	0.51
1:A:986:LYS:HD2	1:A:986:LYS:H	1.74	0.51
1:B:596:SER:OG	1:B:613:GLN:NE2	2.44	0.51
1:B:374:PHE:HA	1:B:436:TRP:HB3	1.93	0.51
2:I:51:ILE:HB	2:I:54:GLY:HA2	1.93	0.51
1:B:642:VAL:HG22	1:B:651:ILE:HG22	1.92	0.50
1:C:985:ASP:OD1	1:C:985:ASP:N	2.42	0.50
3:L:53:ARG:NH2	3:L:58:PRO:O	2.42	0.50
1:A:742:ILE:O	1:A:1000:ARG:NH1	2.44	0.50
1:B:455:LEU:HB2	1:B:493:GLN:HB2	1.94	0.50
1:A:320:VAL:N	1:A:591:SER:O	2.44	0.50
1:A:598:ILE:HB	1:A:609:ALA:HB3	1.94	0.50
1:C:444:LYS:O	1:C:447:GLY:N	2.44	0.50
3:K:19:ARG:NH1	3:K:72:LEU:O	2.44	0.50
2:H:34:MET:HE1	2:H:94:ARG:HG3	1.93	0.50
1:B:36:VAL:HG11	1:B:287:ASP:CB	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:900:MET:SD	1:B:900:MET:N	2.85	0.50
1:A:204:TYR:HB3	1:A:223:LEU:HB3	1.93	0.50
1:C:422:ASN:ND2	1:C:454:ARG:O	2.44	0.50
3:K:53:ARG:NH1	3:K:61:PHE:O	2.43	0.50
1:C:424:LYS:NZ	1:C:425:LEU:O	2.42	0.50
2:I:34:MET:HG3	2:I:94:ARG:HA	1.93	0.50
2:I:52:TRP:O	2:I:71:ARG:NH2	2.44	0.50
1:A:34:ARG:NH2	1:A:218:GLN:O	2.45	0.50
1:A:644:GLN:NE2	1:A:645:THR:O	2.45	0.50
1:B:1047:TYR:HB2	1:B:1067:TYR:HB3	1.94	0.50
1:C:99:ASN:OD1	1:C:190:ARG:NH1	2.45	0.50
1:C:417:LYS:O	1:C:421:TYR:HB2	2.12	0.50
1:C:1030:SER:HA	1:C:1034:LEU:HD13	1.93	0.50
1:A:314:GLN:NE2	1:A:316:SER:O	2.45	0.49
1:A:950:ASP:HA	1:A:953:ASN:HB2	1.93	0.49
2:I:32:TYR:O	2:I:71:ARG:NH2	2.45	0.49
1:C:598:ILE:HB	1:C:609:ALA:HB3	1.94	0.49
1:A:984:LEU:HB3	1:A:988:GLU:HG2	1.94	0.49
1:A:1090:PRO:O	1:B:913:GLN:NE2	2.45	0.49
1:B:815:ARG:HD2	1:B:819:GLU:HB3	1.94	0.49
1:A:69:HIS:H	1:A:78:ARG:HB3	1.77	0.49
1:B:248:TYR:HB2	3:J:95(A):THR:HG21	1.94	0.49
1:A:230:PRO:O	1:C:357:ARG:NH2	2.45	0.49
1:A:781:VAL:HG13	1:A:1029:MET:HE1	1.95	0.49
1:B:702:GLU:HA	1:C:788:ILE:HB	1.95	0.49
1:C:811:LYS:O	1:C:814:LYS:NZ	2.46	0.49
3:K:45:LEU:HD11	3:K:48:PHE:HB3	1.95	0.49
1:A:278:LYS:HD3	1:A:286:THR:H	1.78	0.49
1:A:382:VAL:HG22	1:B:983:ARG:HG2	1.95	0.49
1:A:712:ILE:HG22	1:B:896:ILE:HA	1.95	0.49
1:B:448:ASN:HB2	1:B:497:PHE:HB2	1.93	0.49
1:C:726:ILE:HG12	1:C:945:LEU:HA	1.94	0.49
1:A:563:GLN:H	1:B:40:ASP:HA	1.77	0.49
1:B:124:THR:HG1	7:B:1308:NAG:HO6	1.58	0.49
1:A:108:THR:O	1:A:237:ARG:NH1	2.45	0.49
1:A:584:ILE:HG13	1:B:42:VAL:HA	1.95	0.49
1:A:626:ALA:O	1:A:634:ARG:NH1	2.46	0.49
1:A:1106:GLN:NE2	1:A:1111:GLU:OE2	2.45	0.49
1:B:273:ARG:HH12	1:B:293:LEU:HD11	1.77	0.49
1:C:749:CYS:O	1:C:753:LEU:HB2	2.12	0.49
1:B:503:VAL:HA	1:B:506:GLN:HE22	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1081:ILE:HG23	1:B:1135:ASN:HB3	1.94	0.48
1:C:455:LEU:HD13	1:C:493:GLN:HB2	1.94	0.48
1:C:798:GLY:O	1:C:920:GLN:NE2	2.46	0.48
2:I:12:VAL:HG21	2:I:18:LEU:HB2	1.95	0.48
1:C:915:VAL:O	1:C:919:ASN:ND2	2.46	0.48
1:A:1052:PHE:HB2	1:A:1063:LEU:HB2	1.94	0.48
1:B:1031:GLU:OE1	1:B:1039:ARG:NH2	2.46	0.48
1:C:474:GLN:HA	1:C:488:CYS:HA	1.95	0.48
1:A:41:LYS:HD2	1:C:561:PRO:HD2	1.96	0.48
1:A:202:LYS:HA	1:A:228:ASP:HA	1.94	0.48
1:B:119:ILE:HD11	1:B:175:PHE:HE2	1.79	0.48
1:C:314:GLN:NE2	1:C:316:SER:O	2.45	0.48
3:J:84:ASP:HA	3:J:104:LEU:HD13	1.95	0.48
1:B:710:ASN:OD1	1:B:710:ASN:N	2.46	0.48
1:B:987:VAL:O	1:B:987:VAL:HG12	2.13	0.48
1:B:1118:ASP:OD1	1:B:1118:ASP:N	2.46	0.48
2:H:36:TRP:HE1	2:H:78:LEU:HD22	1.76	0.48
1:A:415:THR:OG1	1:A:416:GLY:N	2.47	0.48
1:A:788:ILE:HB	1:C:702:GLU:HA	1.95	0.48
1:B:172:SER:OG	1:B:173:GLN:N	2.46	0.48
1:C:128:ILE:HB	1:C:170:TYR:HB3	1.95	0.48
2:G:12:VAL:HB	2:G:18:LEU:HD22	1.96	0.48
2:I:38:ARG:NH1	2:I:86:ASP:OD1	2.45	0.48
1:A:321:GLN:NE2	1:A:622:VAL:HB	2.29	0.48
1:A:552:LEU:HB3	1:A:585:LEU:HB3	1.95	0.48
1:A:1053:PRO:O	1:A:1054:GLN:NE2	2.46	0.48
1:B:420:ASP:OD2	1:B:460:ASN:ND2	2.47	0.48
1:B:457:ARG:NH2	1:B:459:SER:OG	2.47	0.48
1:C:287:ASP:N	1:C:287:ASP:OD1	2.46	0.48
1:C:457:ARG:NH2	1:C:459:SER:OG	2.47	0.48
1:C:726:ILE:HA	1:C:1061:VAL:HG12	1.96	0.48
1:B:731:MET:HE2	1:B:774:GLN:HE22	1.79	0.48
1:C:204:TYR:HB3	1:C:223:LEU:HB3	1.95	0.48
1:C:625:HIS:HB3	1:C:628:GLN:HB2	1.96	0.48
1:A:93:ALA:N	1:A:266:TYR:O	2.47	0.47
1:C:743:CYS:HB3	1:C:977:LEU:HD22	1.94	0.47
1:A:710:ASN:OD1	1:A:710:ASN:N	2.47	0.47
1:B:383:SER:HB2	1:C:985:ASP:H	1.79	0.47
3:J:53:ARG:NH1	3:J:73:THR:OG1	2.47	0.47
3:L:92:SER:OG	3:L:93:SER:N	2.44	0.47
1:B:733:LYS:HE2	1:B:863:PRO:HA	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:762:GLN:HG3	1:B:765:ARG:HH21	1.79	0.47
1:C:753:LEU:O	1:C:757:GLY:N	2.47	0.47
2:H:51:ILE:HB	2:H:54:GLY:HA2	1.94	0.47
1:A:112:SER:HA	1:A:132:GLU:HB2	1.96	0.47
1:A:452:LEU:HD23	1:A:492:LEU:HB3	1.96	0.47
1:C:193:VAL:HB	1:C:204:TYR:HB2	1.96	0.47
2:G:100(B):PHE:N	3:J:35:TYR:OH	2.47	0.47
1:B:118:LEU:HB3	1:B:129:LYS:HB2	1.97	0.47
1:B:403:ARG:HG3	1:B:405:ASP:H	1.78	0.47
1:C:146:HIS:HD2	1:C:152:TRP:HB2	1.79	0.47
1:B:427:ASP:OD1	1:B:427:ASP:N	2.47	0.47
1:B:822:LEU:HD11	1:B:1061:VAL:HG21	1.97	0.47
1:A:93:ALA:HB3	1:A:266:TYR:HB2	1.96	0.47
1:A:278:LYS:NZ	1:A:284:THR:O	2.42	0.47
1:A:796:ASP:OD1	1:A:796:ASP:N	2.45	0.47
1:A:951:VAL:O	1:A:955:ASN:ND2	2.44	0.47
1:B:49:HIS:HB3	1:B:277:LEU:HB2	1.97	0.47
1:C:329:PHE:O	1:C:580:GLN:NE2	2.46	0.47
1:C:434:ILE:HB	1:C:511:VAL:HB	1.97	0.47
1:C:673:SER:OG	1:C:674:TYR:N	2.46	0.47
1:C:1091:ARG:NH1	1:C:1118:ASP:O	2.48	0.47
3:K:12:VAL:HB	3:K:14:PRO:HD3	1.95	0.47
2:H:52:TRP:HB2	2:H:55:SER:H	1.79	0.47
1:A:358:ILE:HB	1:A:395:VAL:HB	1.96	0.47
1:B:475:ALA:HB3	1:B:487:ASN:HB3	1.96	0.47
1:B:914:ASN:ND2	1:B:918:GLU:OE2	2.47	0.47
1:B:1054:GLN:HB2	1:B:1061:VAL:HB	1.96	0.47
1:B:328:ARG:NH1	1:B:531:THR:O	2.40	0.47
1:C:334:ASN:ND2	1:C:360:ASN:O	2.46	0.47
3:J:35:TYR:HB2	3:J:86:TYR:HB2	1.96	0.47
3:L:1:SER:OG	3:L:2:TYR:N	2.46	0.47
3:L:6:GLN:NE2	3:L:101:GLY:O	2.48	0.47
1:A:448:ASN:O	1:A:497:PHE:N	2.46	0.47
1:A:808:ASP:OD1	1:A:808:ASP:N	2.46	0.46
1:A:201:PHE:HB3	1:A:229:LEU:HB2	1.97	0.46
1:A:438:SER:HB2	1:A:441:LEU:HB2	1.97	0.46
1:B:1053:PRO:O	1:B:1054:GLN:NE2	2.48	0.46
1:C:16:VAL:HG12	1:C:18:LEU:HB2	1.98	0.46
1:A:528:LYS:NZ	1:A:529:LYS:O	2.48	0.46
1:B:106:PHE:HB2	1:B:117:LEU:HB3	1.96	0.46
1:C:125:ASN:HD22	1:C:171:VAL:HG13	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:448:ASN:OD1	1:C:450:ASN:ND2	2.48	0.46
1:A:32:PHE:H	1:A:216:LEU:HD23	1.80	0.46
1:A:726:ILE:O	1:A:947:LYS:NZ	2.45	0.46
1:A:560:LEU:HB3	1:B:38:TYR:HB3	1.97	0.46
1:B:708:SER:HB2	1:B:711:SER:HB3	1.98	0.46
1:A:319:ARG:HH21	1:A:593:GLY:HA3	1.81	0.46
1:A:559:PHE:HB3	1:A:560:LEU:H	1.43	0.46
1:C:627:ASP:OD1	1:C:627:ASP:N	2.48	0.46
3:J:53:ARG:NH1	3:J:61:PHE:O	2.47	0.46
3:K:23:SER:OG	3:K:24:GLY:N	2.46	0.46
2:H:38:ARG:HG2	2:H:48:VAL:HB	1.97	0.46
1:B:376:THR:HB	1:B:435:ALA:HB3	1.96	0.46
1:C:881:THR:O	1:C:901:GLN:NE2	2.47	0.46
1:B:329:PHE:HB2	1:B:530:SER:HB2	1.98	0.46
2:G:2:VAL:HG21	2:G:94:ARG:HE	1.81	0.46
2:H:52:TRP:O	2:H:71:ARG:NH2	2.49	0.46
1:A:715:PRO:HA	1:A:1072:GLU:HA	1.98	0.46
1:B:276:LEU:O	1:B:288:ALA:CA	2.64	0.46
1:B:727:LEU:HD12	1:B:728:PRO:HD2	1.98	0.46
1:A:376:THR:HB	1:A:435:ALA:HB3	1.98	0.46
1:C:396:TYR:HB2	1:C:514:SER:HB3	1.97	0.46
1:C:905:ARG:NH1	1:C:1049:LEU:O	2.45	0.46
1:A:820:ASP:OD1	1:A:820:ASP:N	2.50	0.45
2:G:89:VAL:HG21	3:J:41:GLN:HE22	1.81	0.45
1:C:654:GLU:HB2	1:C:693:ILE:HA	1.98	0.45
2:H:37:VAL:HA	2:H:47:TRP:HA	1.97	0.45
1:C:438:SER:HB2	1:C:441:LEU:HB2	1.98	0.45
1:C:966:LEU:O	1:C:1000:ARG:NH2	2.48	0.45
1:A:34:ARG:NH1	1:A:221:SER:OG	2.48	0.45
1:A:405:ASP:OD2	1:A:405:ASP:N	2.50	0.45
1:B:417:LYS:HA	1:B:420:ASP:HB2	1.99	0.45
1:C:44:ARG:HB2	1:C:285:ILE:HD11	1.99	0.45
1:C:147:LYS:NZ	3:L:65:SER:O	2.49	0.45
1:C:274:THR:O	1:C:274:THR:OG1	2.35	0.45
1:A:294:ASP:OD1	1:A:297:SER:N	2.48	0.45
1:B:81:ASN:O	1:B:239:GLN:NE2	2.44	0.45
1:B:792:PRO:O	1:B:795:LYS:NZ	2.49	0.45
1:B:1026:ALA:O	1:B:1030:SER:OG	2.34	0.45
1:C:99:ASN:HA	1:C:190:ARG:HH12	1.80	0.45
3:J:19:ARG:HD2	3:J:71:THR:HB	1.98	0.45
1:B:276:LEU:O	1:B:288:ALA:HB3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:52:TRP:CD1	2:I:56:ASN:H	2.35	0.45
1:A:250:THR:HG22	1:A:253:ASP:H	1.82	0.45
1:B:329:PHE:HE2	1:B:528:LYS:HB3	1.82	0.45
2:I:38:ARG:HA	2:I:90:TYR:HA	1.99	0.45
1:A:544:ASN:HD21	1:A:579:PRO:HB3	1.82	0.45
1:B:415:THR:OG1	1:B:416:GLY:N	2.48	0.45
1:C:661:GLU:OE2	1:C:698:SER:OG	2.34	0.45
2:G:36:TRP:N	2:G:49:ALA:O	2.47	0.45
2:I:39:GLN:HB3	2:I:91:TYR:HE1	1.81	0.45
2:I:53:ASP:OD1	2:I:53:ASP:N	2.50	0.45
1:C:346:ARG:NH2	1:C:442:ASP:OD1	2.50	0.45
2:G:82(C):LEU:HD11	2:G:111:VAL:HG11	1.98	0.45
2:H:49:ALA:HB1	2:H:69:ILE:HD13	1.99	0.45
1:A:420:ASP:N	1:A:420:ASP:OD1	2.46	0.44
1:B:206:LYS:HB2	1:B:223:LEU:HA	1.99	0.44
1:B:987:VAL:HG21	1:C:427:ASP:OD2	2.16	0.44
1:C:85:PRO:HA	1:C:237:ARG:HG2	1.98	0.44
1:C:126:VAL:HB	1:C:172:SER:HB3	2.00	0.44
1:C:240:THR:OG1	1:C:241:LEU:N	2.50	0.44
1:C:1105:THR:OG1	1:C:1106:GLN:N	2.50	0.44
1:A:936:ASP:O	1:A:940:SER:N	2.50	0.44
1:B:763:LEU:HD22	1:B:1008:VAL:HG21	1.97	0.44
1:B:836:GLN:NE2	1:B:843:ASP:OD2	2.49	0.44
1:B:375:SER:N	1:B:435:ALA:O	2.50	0.44
1:B:904:TYR:O	1:B:908:GLY:N	2.51	0.44
1:C:201:PHE:H	1:C:229:LEU:H	1.64	0.44
3:K:1:SER:OG	3:K:2:TYR:N	2.49	0.44
1:A:355:ARG:NE	1:A:398:ASP:OD1	2.47	0.44
1:A:1082:CYS:HB3	1:A:1126:CYS:HB3	1.73	0.44
1:B:244:LEU:HD11	1:B:256:SER:HB2	1.99	0.44
3:K:25:ASP:OD1	3:K:25:ASP:N	2.51	0.44
1:B:767:LEU:HA	1:B:770:ILE:HD12	2.00	0.44
1:A:44:ARG:HB2	1:A:285:ILE:HD11	1.99	0.44
1:A:567:ARG:NH1	1:A:571:ASP:O	2.43	0.44
1:B:605:SER:OG	1:B:606:ASN:N	2.50	0.44
3:L:27:LEU:HD11	3:L:70:ILE:HG13	2.00	0.44
1:A:985:ASP:O	1:A:986:LYS:C	2.61	0.44
1:B:327:VAL:HA	1:B:542:ASN:HB3	2.00	0.44
1:C:325:SER:HA	1:C:540:ASN:HB2	2.00	0.44
1:C:353:TRP:O	1:C:466:ARG:NH2	2.45	0.44
1:C:1026:ALA:O	1:C:1030:SER:OG	2.36	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:ASP:OD2	1:A:428:ASP:N	2.51	0.44
1:B:655:HIS:HA	1:B:694:ALA:HB3	1.99	0.44
1:B:1036:GLN:HE21	1:B:1049:LEU:HA	1.82	0.44
2:G:39:GLN:OE1	3:J:41:GLN:NE2	2.50	0.44
1:B:980:ILE:HG23	1:B:984:LEU:HD13	1.99	0.44
1:B:1105:THR:HG23	1:B:1111:GLU:H	1.82	0.44
1:C:327:VAL:HA	1:C:542:ASN:HB3	2.00	0.44
1:B:134:GLN:O	1:B:161:SER:OG	2.36	0.43
1:B:707:TYR:HB2	1:C:883:THR:HG23	1.98	0.43
1:B:707:TYR:HE1	1:C:897:PRO:HA	1.82	0.43
1:A:107:GLY:H	1:A:235:ILE:HG23	1.82	0.43
1:C:867:ASP:N	1:C:867:ASP:OD1	2.48	0.43
3:J:25:ASP:OD1	3:J:25:ASP:N	2.51	0.43
2:H:98:GLY:O	3:L:33:TYR:OH	2.34	0.43
1:A:319:ARG:HB3	1:A:591:SER:HB3	1.99	0.43
1:C:393:THR:HG22	1:C:522:ALA:HA	1.99	0.43
1:C:439:ASN:O	1:C:443:SER:OG	2.35	0.43
2:H:63:VAL:HB	2:H:67:PHE:HE1	1.83	0.43
1:C:93:ALA:O	1:C:266:TYR:N	2.51	0.43
1:C:880:GLY:O	1:C:885:GLY:N	2.46	0.43
2:H:5:VAL:HB	2:H:23:ALA:HB3	2.01	0.43
1:A:300:LYS:HE3	1:A:306:PHE:HA	2.01	0.43
1:B:24:LEU:HA	1:B:25:PRO:HD3	1.84	0.43
1:C:126:VAL:HG23	1:C:174:PRO:HA	2.01	0.43
1:B:359:SER:OG	1:B:360:ASN:N	2.49	0.43
1:B:741:TYR:HA	1:B:856:ASN:HB3	2.01	0.43
1:B:758:SER:O	1:B:762:GLN:NE2	2.52	0.43
3:J:4:LEU:HD11	3:J:89:SER:H	1.84	0.43
3:L:60:ARG:NH2	3:L:80:GLU:OE2	2.44	0.43
1:A:256:SER:OG	1:A:257:GLY:N	2.52	0.43
1:B:960:ASN:O	1:B:964:LYS:HB2	2.18	0.43
1:C:310:LYS:HA	1:C:599:THR:HG23	2.00	0.43
3:L:91:ASP:OD1	3:L:91:ASP:N	2.52	0.43
1:B:962:LEU:HD21	1:B:1007:TYR:HB2	2.01	0.43
1:B:1052:PHE:HB2	1:B:1063:LEU:HB2	2.01	0.43
1:A:43:PHE:HD1	1:C:560:LEU:HG	1.84	0.43
1:A:1093:GLY:O	1:B:904:TYR:OH	2.37	0.43
1:B:359:SER:HA	1:B:524:VAL:HB	2.00	0.43
1:B:742:ILE:O	1:B:1000:ARG:NH1	2.51	0.43
2:I:40:ALA:HA	2:I:88:ALA:HA	2.01	0.43
1:A:320:VAL:N	1:A:591:SER:HB2	2.29	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1052:PHE:HB2	1:C:1063:LEU:HB2	1.99	0.42
4:N:1:NAG:H61	4:N:2:NAG:H82	2.01	0.42
1:A:93:ALA:HA	1:A:191:GLU:HA	2.01	0.42
1:B:49:HIS:HD2	1:B:277:LEU:HD12	1.85	0.42
1:B:598:ILE:N	1:B:609:ALA:O	2.52	0.42
1:C:134:GLN:N	1:C:162:SER:OG	2.49	0.42
3:K:34:TRP:HZ3	3:K:72:LEU:HD13	1.84	0.42
1:A:472:ILE:HA	1:A:491:PRO:HD3	2.00	0.42
1:A:472:ILE:HD13	1:A:482:GLY:H	1.84	0.42
1:A:992:GLN:OE1	1:A:995:ARG:NH1	2.52	0.42
1:B:391:CYS:HA	1:B:525:CYS:HA	2.01	0.42
1:C:29:THR:N	1:C:62:VAL:O	2.46	0.42
2:H:29:PHE:HD2	2:H:73:ASN:HA	1.84	0.42
1:A:631:PRO:HA	1:A:634:ARG:HB2	2.01	0.42
1:A:1116:THR:O	1:A:1120:THR:OG1	2.37	0.42
1:B:77:LYS:HE2	2:G:52(A):TYR:HE1	1.84	0.42
1:B:319:ARG:O	1:B:321:GLN:NE2	2.47	0.42
1:B:818:ILE:O	1:B:822:LEU:HB2	2.19	0.42
1:C:1018:ILE:HD12	1:C:1018:ILE:HA	1.92	0.42
2:G:51:ILE:HG22	2:G:57:LYS:HG2	2.01	0.42
2:I:48:VAL:HG13	2:I:63:VAL:HG21	2.01	0.42
1:A:69:HIS:O	1:A:78:ARG:NE	2.47	0.42
1:B:650:LEU:HD13	1:B:653:ALA:HB3	2.01	0.42
2:I:61:ASP:HA	2:I:64:LYS:HG2	2.01	0.42
1:B:401:VAL:HG22	1:B:509:ARG:HA	2.02	0.42
1:B:529:LYS:NZ	1:B:530:SER:O	2.44	0.42
1:B:712:ILE:HG13	1:C:896:ILE:HA	2.01	0.42
1:C:329:PHE:O	1:C:530:SER:OG	2.37	0.42
1:C:728:PRO:O	1:C:1021:SER:OG	2.36	0.42
3:L:16:GLN:NE2	3:L:17:THR:O	2.53	0.42
1:A:418:ILE:HA	1:A:422:ASN:HB2	2.00	0.42
1:A:1013:ILE:HD11	1:B:1016:ALA:HB2	2.02	0.42
1:A:1080:ALA:HB3	1:A:1129:VAL:HG11	2.01	0.42
1:B:392:PHE:N	1:B:524:VAL:O	2.52	0.42
1:B:725:GLU:HB3	1:B:1062:PHE:HB2	2.02	0.42
1:C:869:MET:HA	1:C:872:GLN:HB2	2.02	0.42
1:A:130:VAL:HG22	1:A:168:PHE:HB3	2.02	0.42
1:A:947:LYS:HE2	1:A:947:LYS:HB3	1.90	0.42
1:B:37:TYR:HA	1:B:223:LEU:HB2	2.02	0.42
1:B:669:GLY:N	1:C:864:LEU:O	2.52	0.42
1:B:731:MET:SD	1:B:731:MET:N	2.93	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:906:PHE:HB3	1:C:911:VAL:HB	2.01	0.42
1:A:534:VAL:HG21	1:A:539:VAL:HG21	2.02	0.42
1:B:667:GLY:HA2	1:C:864:LEU:HA	2.02	0.42
1:C:176:LEU:HD22	1:C:190:ARG:HD2	2.02	0.42
1:C:899:ALA:HA	1:C:902:MET:HE3	2.02	0.42
3:K:65:SER:HA	3:K:70:ILE:HA	2.02	0.42
1:A:415:THR:HA	1:A:419:ALA:HB3	2.00	0.42
1:B:626:ALA:HA	1:B:629:LEU:HB2	2.02	0.42
1:B:721:SER:OG	1:B:1066:THR:OG1	2.38	0.42
1:B:722:VAL:HG12	1:B:930:ALA:HB1	2.01	0.42
1:B:970:PHE:HD2	1:B:996:LEU:HA	1.84	0.42
1:A:318:PHE:N	1:A:318:PHE:CD1	2.87	0.41
1:A:321:GLN:HG2	1:A:623:ALA:H	1.84	0.41
1:A:864:LEU:O	1:C:669:GLY:N	2.53	0.41
1:B:447:GLY:N	1:B:497:PHE:O	2.53	0.41
1:A:100:ILE:HG13	1:A:101:ILE:HG13	2.02	0.41
1:A:669:GLY:N	1:B:864:LEU:O	2.53	0.41
1:C:93:ALA:HA	1:C:191:GLU:HA	2.02	0.41
1:C:146:HIS:HE1	3:L:29:LYS:HE3	1.85	0.41
1:C:250:THR:HG22	1:C:253:ASP:H	1.85	0.41
1:A:767:LEU:HA	1:A:770:ILE:HD12	2.02	0.41
1:A:795:LYS:HA	1:A:795:LYS:HD3	1.96	0.41
1:A:971:GLY:H	1:B:755:GLN:HB3	1.85	0.41
1:B:830:ASP:OD2	1:B:949:GLN:NE2	2.51	0.41
1:C:456:PHE:HB2	1:C:491:PRO:HA	2.02	0.41
2:G:82(C):LEU:HD22	2:G:86:ASP:HB2	2.01	0.41
1:A:303:LEU:HD12	1:A:308:VAL:HG22	2.01	0.41
1:A:320:VAL:HG23	1:A:591:SER:O	2.20	0.41
1:A:627:ASP:O	1:A:634:ARG:NH2	2.52	0.41
1:A:708:SER:OG	1:A:710:ASN:OD1	2.37	0.41
1:A:913:GLN:NE2	1:C:1090:PRO:O	2.54	0.41
1:C:666:ILE:HD11	1:C:672:ALA:HB2	2.01	0.41
1:C:808:ASP:OD2	1:C:810:SER:OG	2.37	0.41
2:G:55:SER:O	4:E:1:NAG:O3	2.37	0.41
1:A:361:CYS:HB2	1:A:524:VAL:HG13	2.03	0.41
1:A:790:LYS:HB2	1:C:704:SER:HA	2.01	0.41
1:B:245:HIS:HD2	3:J:93:SER:HB3	1.85	0.41
1:B:393:THR:HG21	1:B:520:ALA:HB3	2.02	0.41
1:C:815:ARG:HH21	1:C:820:ASP:HA	1.85	0.41
2:I:39:GLN:OE1	2:I:91:TYR:OH	2.39	0.41
1:A:37:TYR:OH	1:A:54:LEU:O	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:633:TRP:HA	1:B:636:TYR:HD2	1.85	0.41
1:C:964:LYS:O	1:C:967:SER:OG	2.36	0.41
1:A:319:ARG:HG2	1:A:591:SER:HB3	2.03	0.41
1:B:134:GLN:N	1:B:162:SER:OG	2.52	0.41
1:B:453:TYR:O	1:B:493:GLN:N	2.52	0.41
3:J:45:LEU:HD21	3:J:48:PHE:HB3	2.02	0.41
3:K:92:SER:OG	3:K:93:SER:N	2.53	0.41
1:A:345:THR:O	1:A:509:ARG:NH2	2.54	0.41
1:B:125:ASN:HD22	1:B:171:VAL:HG13	1.86	0.41
2:I:94:ARG:NH2	2:I:101:ASP:OD2	2.54	0.41
3:L:12:VAL:N	3:L:105:THR:O	2.49	0.41
3:L:25:ASP:OD1	3:L:25:ASP:N	2.54	0.41
3:L:81:ASP:OD1	3:L:81:ASP:N	2.49	0.41
1:A:201:PHE:N	1:A:229:LEU:O	2.54	0.41
1:A:862:PRO:HA	1:A:863:PRO:HD3	1.98	0.41
1:A:914:ASN:ND2	1:A:918:GLU:OE2	2.48	0.41
1:A:1006:THR:HG22	1:A:1010:GLN:HE22	1.84	0.41
1:B:81:ASN:ND2	1:B:240:THR:O	2.54	0.41
1:B:191:GLU:N	1:B:206:LYS:O	2.53	0.41
1:C:203:ILE:HG23	1:C:226:LEU:HB3	2.03	0.41
1:C:275:PHE:HB3	1:C:276:LEU:H	1.74	0.41
1:C:278:LYS:HB3	1:C:285:ILE:HA	2.03	0.41
1:C:475:ALA:HB3	1:C:487:ASN:HB3	2.03	0.41
1:C:761:THR:OG1	1:C:762:GLN:NE2	2.50	0.41
1:A:433:VAL:HG22	1:A:512:VAL:HG12	2.03	0.41
1:B:108:THR:O	1:B:237:ARG:NH2	2.54	0.41
1:B:224:GLU:HA	1:B:225:PRO:HD3	1.95	0.41
1:B:411:ALA:HB3	1:B:414:GLN:HB2	2.01	0.41
1:B:634:ARG:HA	1:B:637:SER:HB3	2.01	0.41
1:B:666:ILE:HD11	1:B:672:ALA:HB2	2.03	0.41
1:B:880:GLY:O	1:B:884:SER:OG	2.35	0.41
1:B:1080:ALA:HB3	1:B:1132:ILE:HG22	2.03	0.41
2:I:56:ASN:HD22	4:F:1:NAG:H83	1.85	0.41
3:K:84:ASP:OD1	3:K:84:ASP:N	2.43	0.41
1:A:41:LYS:HB3	1:C:563:GLN:HB2	2.02	0.40
1:A:96:GLU:CD	1:A:98:SER:H	2.28	0.40
1:A:666:ILE:HD11	1:A:672:ALA:HB2	2.03	0.40
1:B:250:THR:HG22	1:B:253:ASP:H	1.85	0.40
1:B:358:ILE:HG22	1:B:524:VAL:HG21	2.02	0.40
1:B:530:SER:OG	1:B:531:THR:N	2.54	0.40
1:C:374:PHE:HA	1:C:436:TRP:HB3	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1127:ASP:OD1	1:C:1127:ASP:N	2.55	0.40
3:J:61:PHE:O	3:J:73:THR:OG1	2.38	0.40
2:I:6:GLU:O	2:I:105:GLN:NE2	2.55	0.40
3:K:32:GLY:HA2	3:K:89:SER:HB2	2.02	0.40
1:A:318:PHE:HD2	1:A:624:ILE:O	2.05	0.40
1:A:1096:VAL:O	1:A:1103:PHE:N	2.51	0.40
1:B:342:PHE:HE2	1:B:434:ILE:HG21	1.85	0.40
1:C:719:THR:N	1:C:1068:VAL:O	2.50	0.40
3:L:35:TYR:O	3:L:86:TYR:N	2.52	0.40
1:A:286:THR:OG1	1:A:287:ASP:OD1	2.39	0.40
1:B:191:GLU:OE2	1:B:221:SER:OG	2.36	0.40
1:B:438:SER:HB2	1:B:509:ARG:HB2	2.03	0.40
1:B:625:HIS:HB3	1:B:628:GLN:HB2	2.03	0.40
1:C:86:PHE:HE1	1:C:196:ASN:HB2	1.85	0.40
2:I:47:TRP:NE1	2:I:49:ALA:O	2.50	0.40
1:A:40:ASP:HB3	1:A:42:VAL:HG12	2.02	0.40
1:A:555:SER:HB3	1:B:43:PHE:HB2	2.02	0.40
1:A:830:ASP:HA	1:A:849:LEU:HD12	2.04	0.40
2:I:39:GLN:NE2	2:I:40:ALA:O	2.55	0.40
2:H:12:VAL:HG21	2:H:18:LEU:HB2	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	1086/1288 (84%)	997 (92%)	86 (8%)	3 (0%)	36	71
1	B	1089/1288 (84%)	1000 (92%)	87 (8%)	2 (0%)	43	77
1	C	1078/1288 (84%)	1002 (93%)	75 (7%)	1 (0%)	48	83
2	G	117/119 (98%)	101 (86%)	16 (14%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	117/119 (98%)	103 (88%)	14 (12%)	0	100	100
2	I	117/119 (98%)	108 (92%)	9 (8%)	0	100	100
3	J	105/107 (98%)	88 (84%)	17 (16%)	0	100	100
3	K	105/107 (98%)	89 (85%)	16 (15%)	0	100	100
3	L	105/107 (98%)	92 (88%)	13 (12%)	0	100	100
All	All	3919/4542 (86%)	3580 (91%)	333 (8%)	6 (0%)	44	77

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	319	ARG
1	B	288	ALA
1	C	294	ASP
1	A	320	VAL
1	B	290	ASP
1	A	987	VAL

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	951/1114 (85%)	946 (100%)	5 (0%)	81	81
1	B	953/1114 (86%)	952 (100%)	1 (0%)	88	88
1	C	945/1114 (85%)	945 (100%)	0	100	100
2	G	97/97 (100%)	97 (100%)	0	100	100
2	H	97/97 (100%)	95 (98%)	2 (2%)	47	65
2	I	97/97 (100%)	97 (100%)	0	100	100
3	J	89/89 (100%)	89 (100%)	0	100	100
3	K	89/89 (100%)	89 (100%)	0	100	100
3	L	89/89 (100%)	89 (100%)	0	100	100
All	All	3407/3900 (87%)	3399 (100%)	8 (0%)	85	85

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

Mol	Chain	Res	Type
1	A	96	GLU
1	A	318	PHE
1	A	319	ARG
1	A	320	VAL
1	A	986	LYS
1	B	289	VAL
2	H	81	GLN
2	H	105	GLN

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	173	GLN
1	A	314	GLN
1	A	498	GLN
1	A	501	ASN
1	A	506	GLN
1	A	544	ASN
1	A	762	GLN
1	A	853	GLN
1	A	856	ASN
1	B	134	GLN
1	B	334	ASN
1	B	343	ASN
1	B	414	GLN
1	B	439	ASN
1	B	448	ASN
1	B	542	ASN
1	B	613	GLN
1	B	625	HIS
1	B	644	GLN
1	B	804	GLN
1	B	919	ASN
1	B	1023	ASN
1	C	450	ASN
1	C	556	ASN
1	C	607	GLN
1	C	856	ASN
1	C	935	GLN
1	C	955	ASN
1	C	969	ASN

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Mol	Chain	Res	Type
1	C	1106	GLN
3	J	41	GLN
2	I	73	ASN
3	K	6	GLN
2	H	73	ASN
2	H	105	GLN
3	L	41	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

33 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	$\# Z > 2$
5	NAG	D	1	5,1	14,14,15	1.19	2 (14%)	17,19,21	0.82	1 (5%)
5	NAG	D	2	5	14,14,15	1.35	2 (14%)	17,19,21	1.75	3 (17%)
4	NAG	E	1	4,1	14,14,15	1.00	1 (7%)	17,19,21	1.33	3 (17%)
4	NAG	E	2	4	14,14,15	1.04	1 (7%)	17,19,21	0.93	1 (5%)
4	BMA	E	3	4	11,11,12	1.80	4 (36%)	15,15,17	1.24	1 (6%)
4	MAN	E	4	4	11,11,12	1.51	3 (27%)	15,15,17	1.59	1 (6%)
4	NAG	E	5	4	14,14,15	1.49	2 (14%)	17,19,21	1.61	3 (17%)
4	NAG	E	6	4	14,14,15	1.12	1 (7%)	17,19,21	0.94	1 (5%)
4	BMA	E	7	4	11,11,12	1.52	2 (18%)	15,15,17	1.12	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FUC	E	8	4	10,10,11	1.48	3 (30%)	14,14,16	1.02	1 (7%)
4	NAG	F	1	4,1	14,14,15	1.51	2 (14%)	17,19,21	1.34	2 (11%)
4	NAG	F	2	4	14,14,15	0.98	1 (7%)	17,19,21	0.86	1 (5%)
4	BMA	F	3	4	11,11,12	1.59	3 (27%)	15,15,17	1.30	1 (6%)
4	MAN	F	4	4	11,11,12	1.28	1 (9%)	15,15,17	1.71	1 (6%)
4	NAG	F	5	4	14,14,15	1.48	2 (14%)	17,19,21	1.65	3 (17%)
4	NAG	F	6	4	14,14,15	1.09	1 (7%)	17,19,21	1.02	1 (5%)
4	BMA	F	7	4	11,11,12	1.51	2 (18%)	15,15,17	1.18	2 (13%)
4	FUC	F	8	4	10,10,11	1.65	2 (20%)	14,14,16	1.42	2 (14%)
5	NAG	M	1	5,1	14,14,15	1.01	1 (7%)	17,19,21	0.92	1 (5%)
5	NAG	M	2	5	14,14,15	1.46	2 (14%)	17,19,21	1.70	3 (17%)
4	NAG	N	1	4,1	14,14,15	0.83	1 (7%)	17,19,21	1.48	4 (23%)
4	NAG	N	2	4	14,14,15	0.93	1 (7%)	17,19,21	0.97	1 (5%)
4	BMA	N	3	4	11,11,12	1.69	3 (27%)	15,15,17	1.37	3 (20%)
4	MAN	N	4	4	11,11,12	1.46	2 (18%)	15,15,17	1.83	2 (13%)
4	NAG	N	5	4	14,14,15	1.34	2 (14%)	17,19,21	1.67	3 (17%)
4	NAG	N	6	4	14,14,15	0.98	1 (7%)	17,19,21	1.02	1 (5%)
4	BMA	N	7	4	11,11,12	1.41	1 (9%)	15,15,17	1.13	2 (13%)
4	FUC	N	8	4	10,10,11	1.59	2 (20%)	14,14,16	1.52	2 (14%)
5	NAG	O	1	5,1	14,14,15	1.03	1 (7%)	17,19,21	1.08	1 (5%)
5	NAG	O	2	5	14,14,15	1.36	2 (14%)	17,19,21	1.71	3 (17%)
6	NAG	Y	1	6,1	14,14,15	0.89	1 (7%)	17,19,21	1.47	2 (11%)
6	NAG	Y	2	6	14,14,15	1.27	2 (14%)	17,19,21	2.00	3 (17%)
6	BMA	Y	3	6	11,11,12	1.35	2 (18%)	15,15,17	1.14	1 (6%)

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	D	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	D	2	5	-	4/6/23/26	0/1/1/1
4	NAG	E	1	4,1	-	4/6/23/26	0/1/1/1
4	NAG	E	2	4	-	2/6/23/26	0/1/1/1
4	BMA	E	3	4	-	1/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MAN	E	4	4	-	2/2/19/22	0/1/1/1
4	NAG	E	5	4	-	3/6/23/26	0/1/1/1
4	NAG	E	6	4	-	1/6/23/26	0/1/1/1
4	BMA	E	7	4	-	0/2/19/22	0/1/1/1
4	FUC	E	8	4	-	-	0/1/1/1
4	NAG	F	1	4,1	-	4/6/23/26	0/1/1/1
4	NAG	F	2	4	-	1/6/23/26	0/1/1/1
4	BMA	F	3	4	-	0/2/19/22	0/1/1/1
4	MAN	F	4	4	-	1/2/19/22	0/1/1/1
4	NAG	F	5	4	-	4/6/23/26	0/1/1/1
4	NAG	F	6	4	-	2/6/23/26	0/1/1/1
4	BMA	F	7	4	-	0/2/19/22	0/1/1/1
4	FUC	F	8	4	-	-	0/1/1/1
5	NAG	M	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	M	2	5	-	4/6/23/26	0/1/1/1
4	NAG	N	1	4,1	-	3/6/23/26	0/1/1/1
4	NAG	N	2	4	-	2/6/23/26	0/1/1/1
4	BMA	N	3	4	-	0/2/19/22	0/1/1/1
4	MAN	N	4	4	-	1/2/19/22	0/1/1/1
4	NAG	N	5	4	-	4/6/23/26	0/1/1/1
4	NAG	N	6	4	-	2/6/23/26	0/1/1/1
4	BMA	N	7	4	-	0/2/19/22	0/1/1/1
4	FUC	N	8	4	-	-	0/1/1/1
5	NAG	O	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	O	2	5	-	2/6/23/26	0/1/1/1
6	NAG	Y	1	6,1	-	4/6/23/26	0/1/1/1
6	NAG	Y	2	6	-	3/6/23/26	0/1/1/1
6	BMA	Y	3	6	-	2/2/19/22	0/1/1/1

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	5	NAG	O5-C1	4.50	1.51	1.43
4	F	1	NAG	O5-C1	4.16	1.50	1.43
4	N	5	NAG	O5-C1	4.14	1.50	1.43
4	E	5	NAG	O5-C1	4.09	1.50	1.43
4	F	8	FUC	O5-C5	3.89	1.51	1.43
5	M	2	NAG	C1-C2	3.76	1.57	1.52
4	N	8	FUC	O5-C5	3.73	1.51	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	6	NAG	O5-C1	3.68	1.49	1.43
5	M	2	NAG	O5-C1	3.62	1.49	1.43
4	E	6	NAG	O5-C1	3.56	1.49	1.43
5	D	1	NAG	O5-C1	3.55	1.49	1.43
4	F	1	NAG	C1-C2	3.54	1.57	1.52
6	Y	2	NAG	O5-C1	3.45	1.49	1.43
5	O	2	NAG	C1-C2	3.39	1.57	1.52
5	O	2	NAG	O5-C1	3.37	1.49	1.43
5	D	2	NAG	O5-C1	3.37	1.49	1.43
5	D	2	NAG	C1-C2	3.34	1.56	1.52
5	O	1	NAG	O5-C1	3.34	1.49	1.43
4	E	5	NAG	C1-C2	3.34	1.56	1.52
4	N	3	BMA	O5-C5	3.28	1.49	1.43
4	N	4	MAN	O5-C5	3.27	1.49	1.43
4	F	3	BMA	O5-C5	3.26	1.49	1.43
4	N	6	NAG	O5-C1	3.15	1.49	1.43
4	E	3	BMA	C2-C3	3.14	1.57	1.52
5	M	1	NAG	O5-C1	3.13	1.48	1.43
4	E	2	NAG	O5-C1	3.12	1.48	1.43
4	F	2	NAG	O5-C1	3.05	1.48	1.43
4	F	7	BMA	C2-C3	3.01	1.57	1.52
4	E	3	BMA	O5-C5	2.95	1.49	1.43
4	N	3	BMA	C2-C3	2.93	1.57	1.52
4	E	4	MAN	O5-C5	2.91	1.49	1.43
6	Y	2	NAG	C1-C2	2.90	1.56	1.52
4	E	7	BMA	C2-C3	2.88	1.56	1.52
4	N	2	NAG	O5-C1	2.86	1.48	1.43
4	F	3	BMA	C2-C3	2.84	1.56	1.52
4	F	5	NAG	C1-C2	2.76	1.56	1.52
4	E	8	FUC	C2-C3	2.66	1.56	1.52
4	F	4	MAN	O5-C5	2.66	1.48	1.43
4	N	7	BMA	C2-C3	2.65	1.56	1.52
4	F	8	FUC	O5-C1	2.62	1.48	1.43
4	E	1	NAG	C1-C2	2.62	1.55	1.52
6	Y	1	NAG	O5-C1	2.61	1.48	1.43
4	E	3	BMA	C4-C5	2.57	1.58	1.53
4	N	4	MAN	O5-C1	2.52	1.47	1.43
6	Y	3	BMA	C2-C3	2.48	1.56	1.52
4	N	1	NAG	C1-C2	2.47	1.55	1.52
4	E	8	FUC	O5-C5	2.41	1.48	1.43
4	N	5	NAG	C1-C2	2.31	1.55	1.52
4	E	7	BMA	C1-C2	2.23	1.57	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	4	MAN	C1-C2	2.20	1.57	1.52
4	E	8	FUC	C4-C3	2.19	1.58	1.52
4	N	8	FUC	O5-C1	2.17	1.47	1.43
5	D	1	NAG	C1-C2	2.16	1.55	1.52
4	E	3	BMA	C4-C3	2.13	1.57	1.52
6	Y	3	BMA	C4-C3	2.13	1.57	1.52
4	E	4	MAN	C2-C3	2.10	1.55	1.52
4	F	3	BMA	C4-C3	2.10	1.57	1.52
4	F	7	BMA	C4-C3	2.07	1.57	1.52
4	N	3	BMA	C4-C5	2.01	1.57	1.53

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	N	4	MAN	C1-O5-C5	6.29	120.61	112.19
4	F	4	MAN	C1-O5-C5	5.70	119.83	112.19
4	E	4	MAN	C1-O5-C5	5.34	119.34	112.19
4	N	5	NAG	C2-N2-C7	5.14	129.78	122.90
6	Y	2	NAG	C2-N2-C7	5.14	129.78	122.90
5	O	2	NAG	C2-N2-C7	5.12	129.75	122.90
4	F	5	NAG	C2-N2-C7	5.07	129.69	122.90
5	M	2	NAG	C2-N2-C7	5.05	129.67	122.90
4	E	5	NAG	C2-N2-C7	5.02	129.62	122.90
5	D	2	NAG	C2-N2-C7	5.00	129.60	122.90
6	Y	2	NAG	C1-O5-C5	4.97	118.85	112.19
6	Y	1	NAG	C1-O5-C5	4.52	118.24	112.19
4	F	8	FUC	C1-O5-C5	4.14	122.73	112.97
4	N	8	FUC	C1-O5-C5	4.08	122.58	112.97
5	D	2	NAG	C1-O5-C5	3.96	117.50	112.19
4	F	6	NAG	C1-O5-C5	3.92	117.44	112.19
4	N	6	NAG	C1-O5-C5	3.91	117.42	112.19
4	F	3	BMA	C1-O5-C5	3.83	117.32	112.19
4	N	1	NAG	C1-O5-C5	3.77	117.23	112.19
4	E	3	BMA	C1-O5-C5	3.65	117.08	112.19
4	E	6	NAG	C1-O5-C5	3.57	116.97	112.19
4	F	1	NAG	C2-N2-C7	3.54	127.65	122.90
4	N	3	BMA	C1-O5-C5	3.53	116.92	112.19
5	M	2	NAG	C1-O5-C5	3.46	116.82	112.19
4	N	2	NAG	C1-O5-C5	3.43	116.79	112.19
5	O	1	NAG	C1-O5-C5	3.36	116.69	112.19
5	O	2	NAG	C1-O5-C5	3.34	116.67	112.19
4	E	2	NAG	C1-O5-C5	3.34	116.66	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	N	5	NAG	C1-O5-C5	3.33	116.65	112.19
6	Y	3	BMA	C1-O5-C5	3.21	116.49	112.19
4	F	5	NAG	C1-O5-C5	3.17	116.43	112.19
4	N	8	FUC	O5-C5-C4	3.07	115.08	109.55
4	F	7	BMA	C1-O5-C5	3.02	116.24	112.19
4	N	7	BMA	C1-O5-C5	3.02	116.24	112.19
4	F	2	NAG	C1-O5-C5	3.00	116.21	112.19
4	N	1	NAG	C2-N2-C7	2.89	126.78	122.90
4	E	5	NAG	C1-O5-C5	2.88	116.05	112.19
6	Y	1	NAG	C2-N2-C7	2.87	126.75	122.90
4	E	1	NAG	C2-N2-C7	2.87	126.75	122.90
6	Y	2	NAG	C1-C2-N2	2.85	114.92	110.43
4	E	7	BMA	C1-O5-C5	2.84	115.99	112.19
5	M	2	NAG	C1-C2-N2	2.82	114.88	110.43
5	O	2	NAG	C1-C2-N2	2.80	114.84	110.43
4	E	1	NAG	C4-C3-C2	2.75	115.04	111.02
5	M	1	NAG	C1-O5-C5	2.74	115.86	112.19
4	F	1	NAG	C1-O5-C5	2.69	115.78	112.19
4	F	8	FUC	O5-C5-C4	2.66	114.34	109.55
5	D	2	NAG	C1-C2-N2	2.65	114.61	110.43
4	F	5	NAG	C1-C2-N2	2.64	114.59	110.43
4	E	1	NAG	C1-O5-C5	2.57	115.63	112.19
4	E	5	NAG	C1-C2-N2	2.49	114.36	110.43
5	D	1	NAG	C1-O5-C5	2.47	115.50	112.19
4	N	3	BMA	C1-C2-C3	-2.46	106.06	109.64
4	N	1	NAG	C4-C3-C2	2.46	114.62	111.02
4	N	3	BMA	O3-C3-C2	2.43	115.02	110.05
4	N	5	NAG	C1-C2-N2	2.42	114.25	110.43
4	E	8	FUC	C1-O5-C5	2.38	118.59	112.97
4	N	4	MAN	O2-C2-C3	-2.12	105.77	110.15
4	E	7	BMA	C2-C3-C4	2.10	114.55	110.86
4	N	1	NAG	C3-C4-C5	2.07	113.98	110.23
4	F	7	BMA	C2-C3-C4	2.07	114.50	110.86
4	N	7	BMA	C2-C3-C4	2.07	114.50	110.86

There are no chirality outliers.

All (61) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	N	5	NAG	O5-C5-C6-O6
4	F	5	NAG	C4-C5-C6-O6
4	F	1	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	D	2	NAG	C4-C5-C6-O6
6	Y	1	NAG	O5-C5-C6-O6
5	M	2	NAG	C4-C5-C6-O6
5	O	1	NAG	O5-C5-C6-O6
4	F	5	NAG	O5-C5-C6-O6
4	N	5	NAG	C4-C5-C6-O6
4	F	6	NAG	O5-C5-C6-O6
4	N	6	NAG	O5-C5-C6-O6
5	D	2	NAG	O5-C5-C6-O6
5	O	1	NAG	C4-C5-C6-O6
4	F	6	NAG	C4-C5-C6-O6
4	N	6	NAG	C4-C5-C6-O6
6	Y	1	NAG	C4-C5-C6-O6
4	F	1	NAG	O5-C5-C6-O6
5	D	1	NAG	O5-C5-C6-O6
5	M	2	NAG	O5-C5-C6-O6
4	E	1	NAG	C8-C7-N2-C2
4	E	1	NAG	O7-C7-N2-C2
4	F	1	NAG	C8-C7-N2-C2
4	F	1	NAG	O7-C7-N2-C2
4	N	1	NAG	C8-C7-N2-C2
4	N	1	NAG	O7-C7-N2-C2
6	Y	1	NAG	C8-C7-N2-C2
6	Y	1	NAG	O7-C7-N2-C2
4	N	2	NAG	O5-C5-C6-O6
4	E	4	MAN	O5-C5-C6-O6
4	F	4	MAN	O5-C5-C6-O6
4	E	2	NAG	O5-C5-C6-O6
4	N	4	MAN	O5-C5-C6-O6
6	Y	3	BMA	C4-C5-C6-O6
4	F	2	NAG	O5-C5-C6-O6
4	E	6	NAG	O5-C5-C6-O6
4	E	5	NAG	O5-C5-C6-O6
5	M	1	NAG	O5-C5-C6-O6
5	D	1	NAG	C4-C5-C6-O6
6	Y	2	NAG	O5-C5-C6-O6
6	Y	3	BMA	O5-C5-C6-O6
4	E	1	NAG	O5-C5-C6-O6
4	E	1	NAG	C4-C5-C6-O6
5	M	2	NAG	C1-C2-N2-C7
6	Y	2	NAG	C1-C2-N2-C7
4	N	2	NAG	C4-C5-C6-O6

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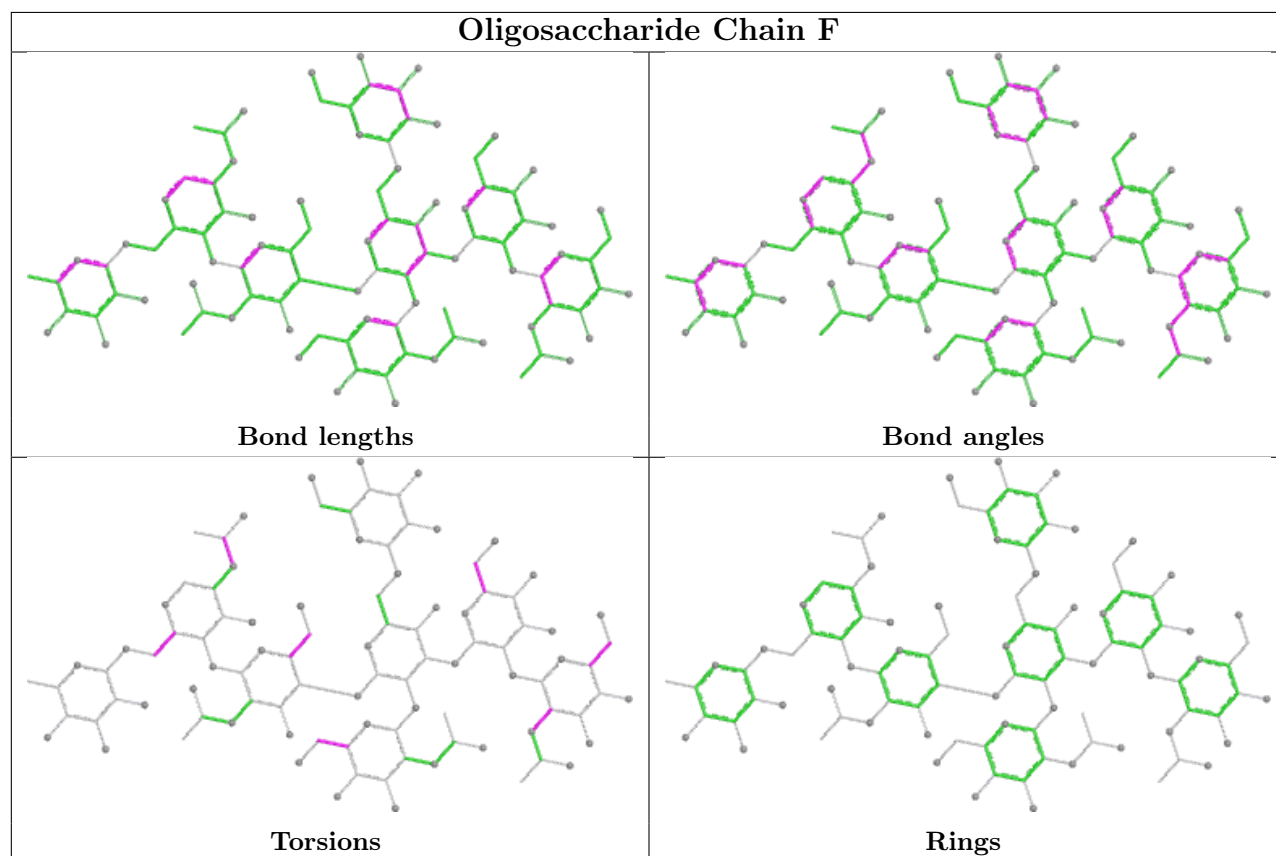
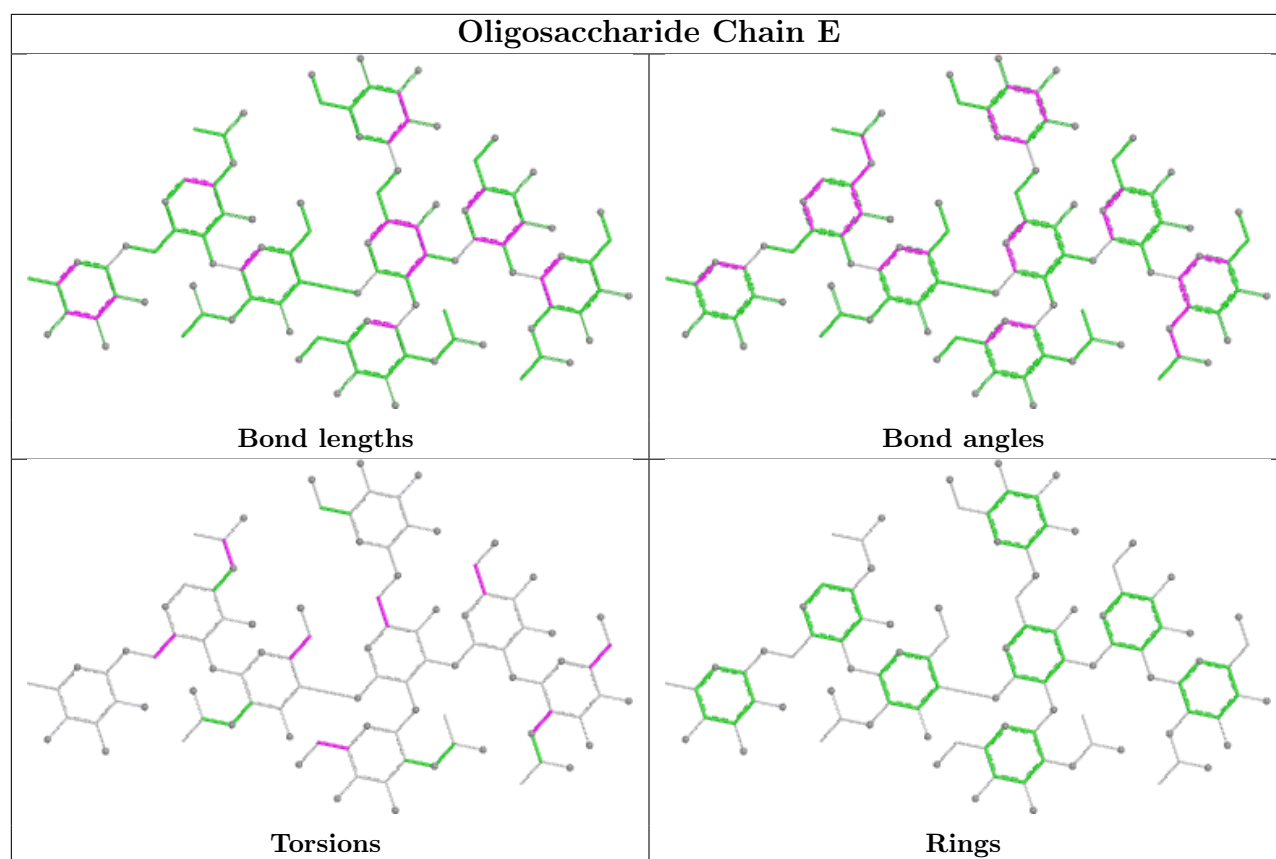
Mol	Chain	Res	Type	Atoms
4	F	5	NAG	C3-C2-N2-C7
4	N	5	NAG	C3-C2-N2-C7
5	M	2	NAG	C3-C2-N2-C7
5	O	2	NAG	C3-C2-N2-C7
6	Y	2	NAG	C3-C2-N2-C7
4	E	3	BMA	C4-C5-C6-O6
4	E	2	NAG	C4-C5-C6-O6
4	E	4	MAN	C4-C5-C6-O6
4	E	5	NAG	C1-C2-N2-C7
4	F	5	NAG	C1-C2-N2-C7
4	N	5	NAG	C1-C2-N2-C7
5	D	2	NAG	C1-C2-N2-C7
5	O	2	NAG	C1-C2-N2-C7
4	E	5	NAG	C3-C2-N2-C7
5	D	2	NAG	C3-C2-N2-C7
4	N	1	NAG	C4-C5-C6-O6

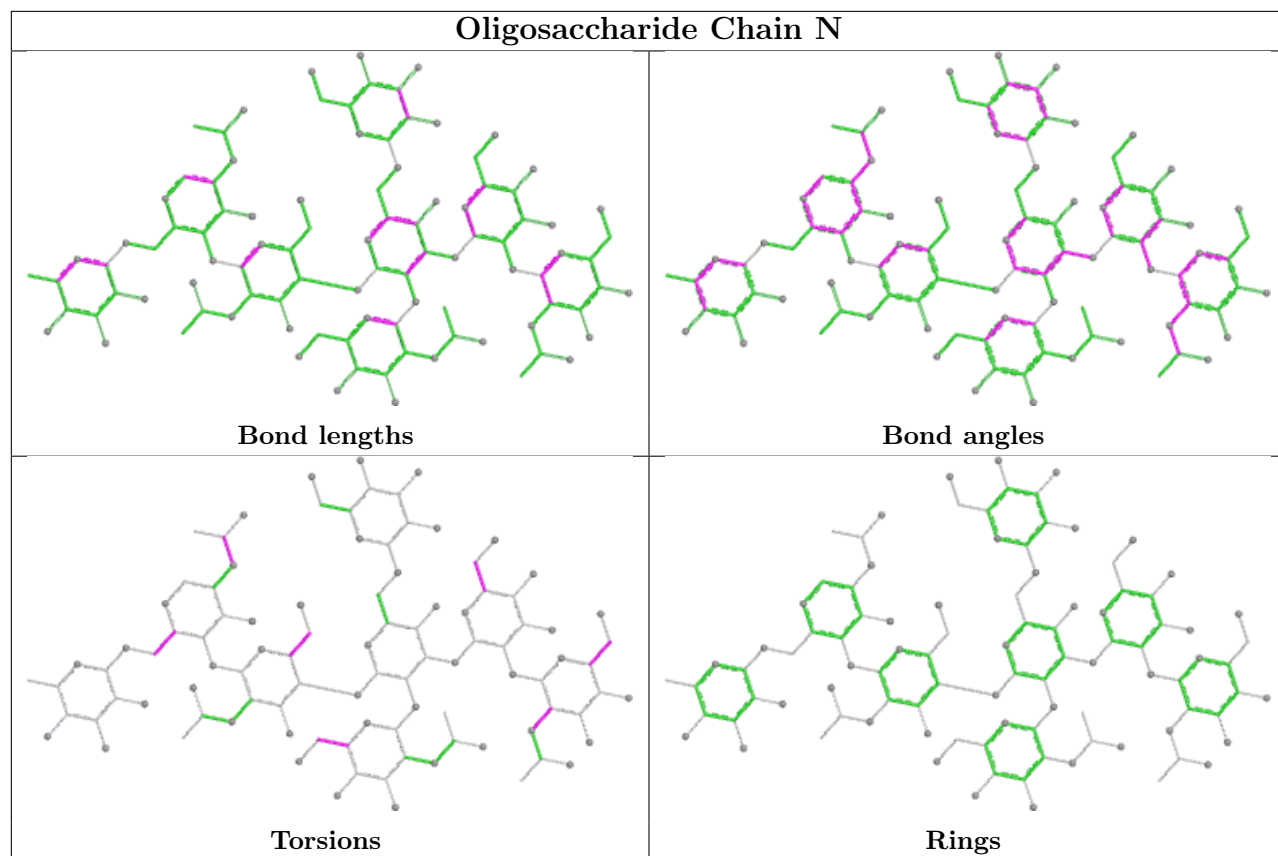
There are no ring outliers.

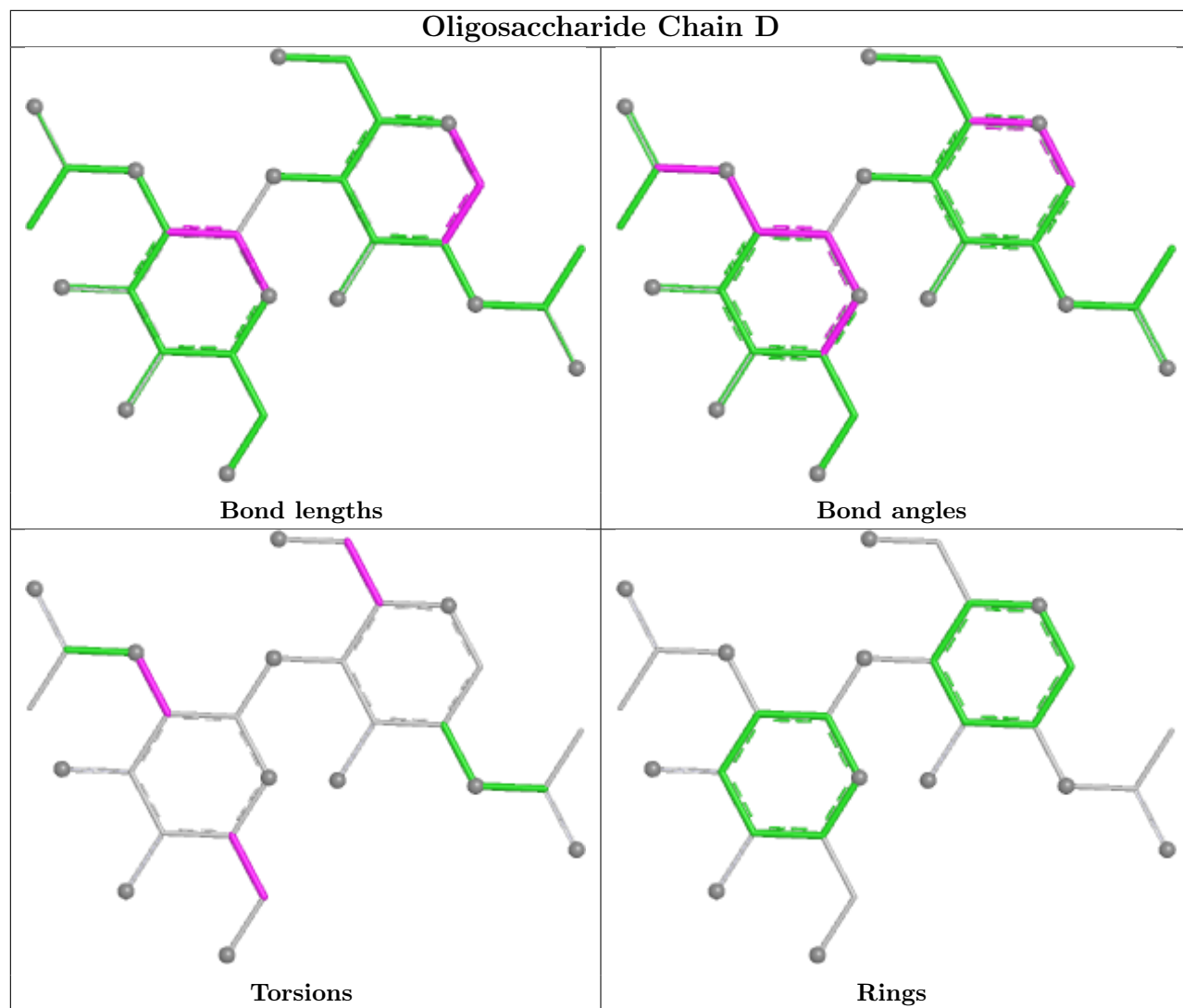
4 monomers are involved in 3 short contacts:

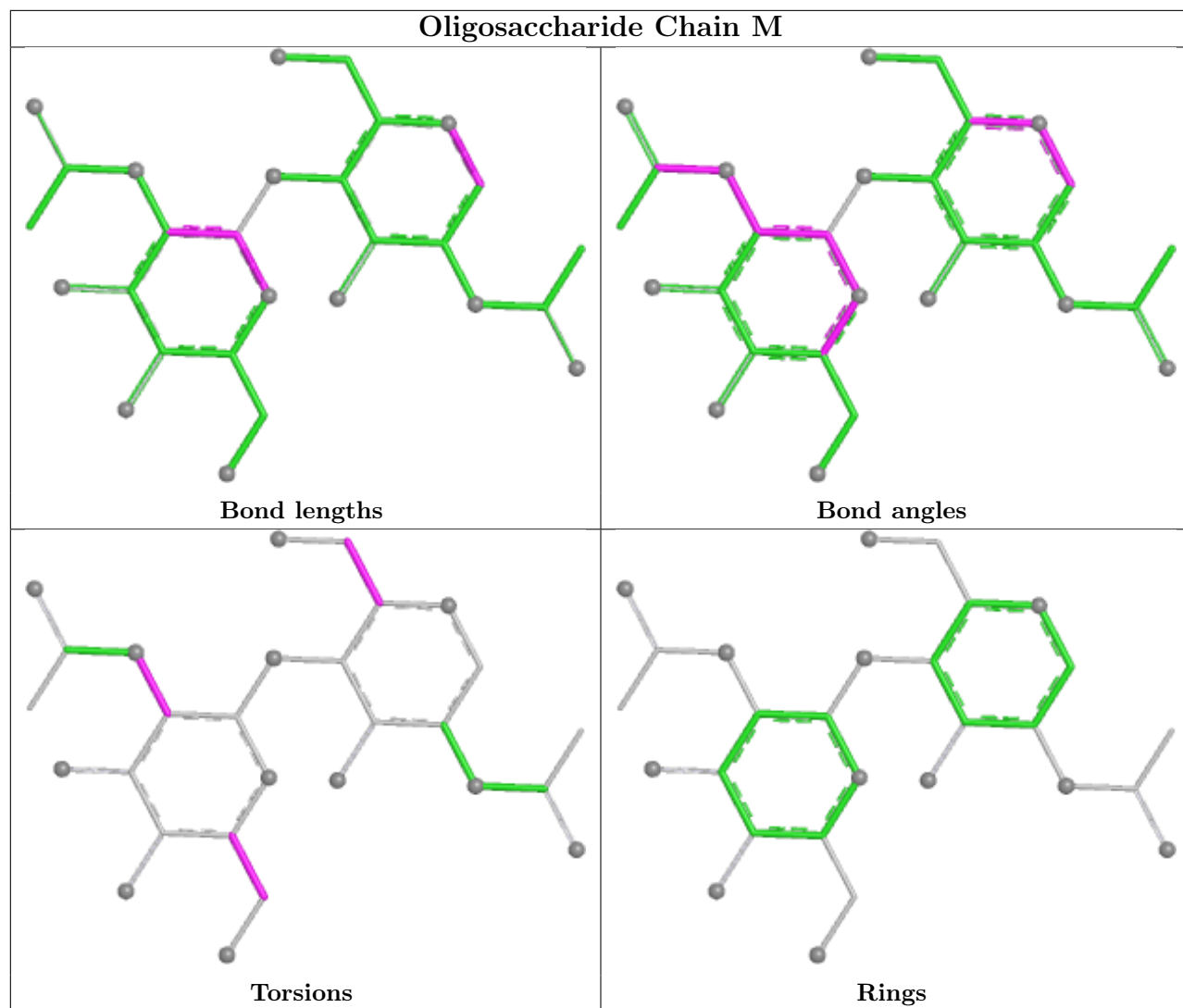
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	1	NAG	1	0
4	N	2	NAG	1	0
4	E	1	NAG	1	0
4	N	1	NAG	1	0

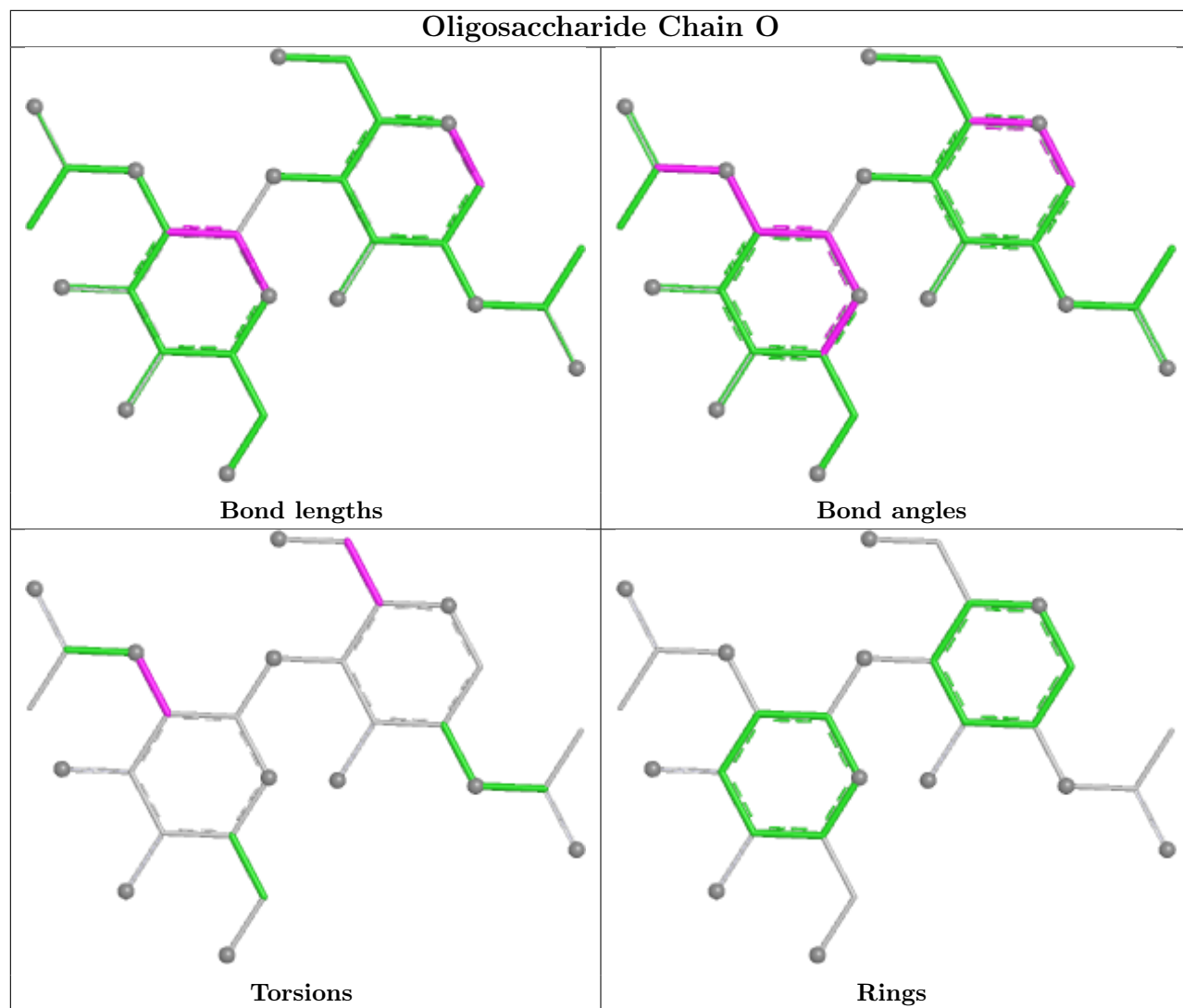
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

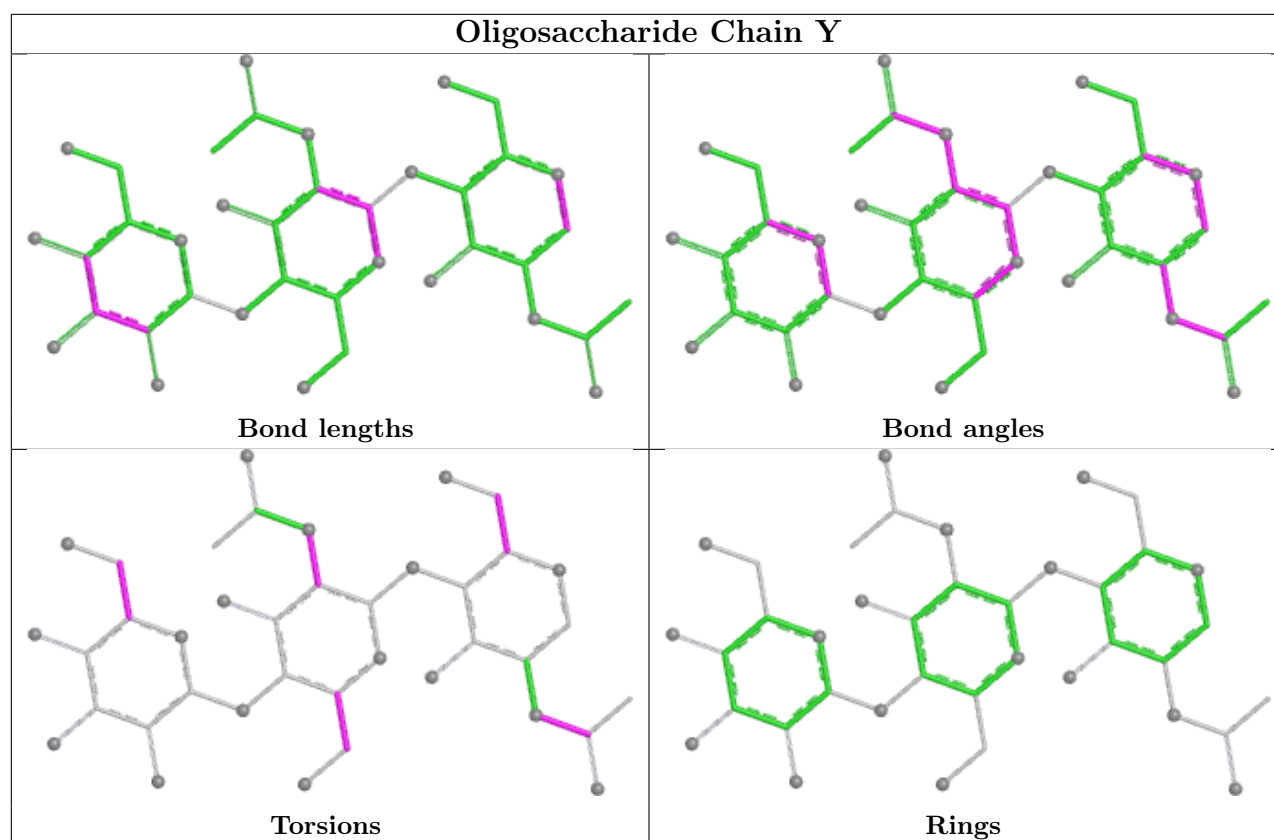












5.6 Ligand geometry [i](#)

28 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
7	NAG	C	1306	1	14,14,15	1.04	1 (7%)	17,19,21	0.87	1 (5%)
7	NAG	B	1304	1	14,14,15	1.20	2 (14%)	17,19,21	1.62	3 (17%)
7	NAG	A	1308	1	14,14,15	1.14	2 (14%)	17,19,21	1.73	3 (17%)
7	NAG	C	1301	-	14,14,15	1.08	1 (7%)	17,19,21	0.91	1 (5%)
7	NAG	C	1308	1	14,14,15	1.12	2 (14%)	17,19,21	1.68	3 (17%)
7	NAG	A	1303	1	14,14,15	1.04	1 (7%)	17,19,21	1.00	1 (5%)
7	NAG	B	1305	1	14,14,15	1.04	1 (7%)	17,19,21	0.87	1 (5%)
7	NAG	B	1306	1	14,14,15	1.11	1 (7%)	17,19,21	0.98	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NAG	B	1309	1	14,14,15	1.04	1 (7%)	17,19,21	1.60	3 (17%)
7	NAG	B	1308	1	14,14,15	1.63	2 (14%)	17,19,21	1.55	2 (11%)
7	NAG	C	1305	1	14,14,15	1.08	1 (7%)	17,19,21	0.82	1 (5%)
7	NAG	A	1307	1	14,14,15	1.53	2 (14%)	17,19,21	1.55	2 (11%)
7	NAG	A	1306	1	14,14,15	1.35	2 (14%)	17,19,21	0.94	1 (5%)
7	NAG	C	1304	1	14,14,15	0.95	1 (7%)	17,19,21	0.99	1 (5%)
7	NAG	A	1302	-	14,14,15	1.04	1 (7%)	17,19,21	0.89	1 (5%)
7	NAG	B	1310	1	14,14,15	1.29	2 (14%)	17,19,21	0.91	1 (5%)
7	NAG	A	1309	1	14,14,15	1.33	2 (14%)	17,19,21	1.16	1 (5%)
7	NAG	A	1301	1	14,14,15	1.22	1 (7%)	17,19,21	1.14	1 (5%)
7	NAG	B	1303	1	14,14,15	1.09	2 (14%)	17,19,21	1.73	3 (17%)
7	NAG	B	1307	-	14,14,15	1.11	1 (7%)	17,19,21	0.87	1 (5%)
7	NAG	C	1302	1	14,14,15	1.20	1 (7%)	17,19,21	0.70	1 (5%)
7	NAG	B	1301	-	14,14,15	1.35	2 (14%)	17,19,21	0.81	1 (5%)
7	NAG	B	1302	1	14,14,15	1.15	1 (7%)	17,19,21	1.08	1 (5%)
7	NAG	C	1303	1	14,14,15	1.04	2 (14%)	17,19,21	1.01	1 (5%)
7	NAG	A	1305	1	14,14,15	1.13	1 (7%)	17,19,21	0.77	1 (5%)
7	NAG	C	1307	1	14,14,15	1.08	2 (14%)	17,19,21	1.02	1 (5%)
7	NAG	A	1304	1	14,14,15	1.06	1 (7%)	17,19,21	0.94	1 (5%)
7	NAG	C	1309	1	14,14,15	1.18	2 (14%)	17,19,21	0.89	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	C	1306	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1304	1	-	3/6/23/26	0/1/1/1
7	NAG	A	1308	1	-	4/6/23/26	0/1/1/1
7	NAG	C	1301	-	-	2/6/23/26	0/1/1/1
7	NAG	C	1308	1	-	4/6/23/26	0/1/1/1
7	NAG	A	1303	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1305	1	-	0/6/23/26	0/1/1/1
7	NAG	B	1306	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1309	1	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	B	1308	1	-	4/6/23/26	0/1/1/1
7	NAG	C	1305	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1307	1	-	4/6/23/26	0/1/1/1
7	NAG	A	1306	1	-	2/6/23/26	0/1/1/1
7	NAG	C	1304	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1302	-	-	2/6/23/26	0/1/1/1
7	NAG	B	1310	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1309	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1301	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1303	1	-	4/6/23/26	0/1/1/1
7	NAG	B	1307	-	-	2/6/23/26	0/1/1/1
7	NAG	C	1302	1	-	2/6/23/26	0/1/1/1
7	NAG	B	1301	-	-	2/6/23/26	0/1/1/1
7	NAG	B	1302	1	-	2/6/23/26	0/1/1/1
7	NAG	C	1303	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1305	1	-	2/6/23/26	0/1/1/1
7	NAG	C	1307	1	-	2/6/23/26	0/1/1/1
7	NAG	A	1304	1	-	0/6/23/26	0/1/1/1
7	NAG	C	1309	1	-	2/6/23/26	0/1/1/1

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	1308	NAG	O5-C1	5.13	1.52	1.43
7	A	1307	NAG	O5-C1	5.00	1.52	1.43
7	B	1301	NAG	O5-C1	4.45	1.51	1.43
7	A	1301	NAG	O5-C1	4.14	1.50	1.43
7	B	1310	NAG	O5-C1	4.02	1.50	1.43
7	C	1302	NAG	O5-C1	3.87	1.50	1.43
7	B	1302	NAG	O5-C1	3.72	1.49	1.43
7	A	1309	NAG	C1-C2	3.65	1.57	1.52
7	A	1305	NAG	O5-C1	3.63	1.49	1.43
7	A	1306	NAG	O5-C1	3.55	1.49	1.43
7	C	1301	NAG	O5-C1	3.54	1.49	1.43
7	C	1309	NAG	O5-C1	3.52	1.49	1.43
7	B	1306	NAG	O5-C1	3.50	1.49	1.43
7	B	1304	NAG	O5-C1	3.49	1.49	1.43
7	B	1307	NAG	O5-C1	3.46	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	1305	NAG	O5-C1	3.42	1.49	1.43
7	A	1304	NAG	O5-C1	3.40	1.49	1.43
7	A	1306	NAG	C1-C2	3.34	1.56	1.52
7	B	1305	NAG	O5-C1	3.33	1.49	1.43
7	A	1302	NAG	O5-C1	3.32	1.49	1.43
7	B	1309	NAG	O5-C1	3.26	1.49	1.43
7	C	1307	NAG	O5-C1	3.23	1.49	1.43
7	A	1309	NAG	O5-C1	3.17	1.49	1.43
7	A	1303	NAG	O5-C1	3.16	1.49	1.43
7	C	1303	NAG	O5-C1	3.11	1.48	1.43
7	C	1306	NAG	O5-C1	3.10	1.48	1.43
7	A	1308	NAG	O5-C1	3.09	1.48	1.43
7	C	1308	NAG	O5-C1	3.04	1.48	1.43
7	B	1303	NAG	O5-C1	2.99	1.48	1.43
7	B	1308	NAG	C1-C2	2.91	1.56	1.52
7	C	1304	NAG	O5-C1	2.73	1.48	1.43
7	A	1308	NAG	C1-C2	2.44	1.55	1.52
7	C	1308	NAG	C1-C2	2.39	1.55	1.52
7	C	1309	NAG	C1-C2	2.38	1.55	1.52
7	A	1307	NAG	C1-C2	2.35	1.55	1.52
7	B	1303	NAG	C1-C2	2.32	1.55	1.52
7	B	1310	NAG	C1-C2	2.28	1.55	1.52
7	B	1304	NAG	C1-C2	2.26	1.55	1.52
7	C	1307	NAG	C1-C2	2.09	1.55	1.52
7	B	1301	NAG	C1-C2	2.02	1.55	1.52
7	C	1303	NAG	C1-C2	2.01	1.55	1.52

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	1308	NAG	C2-N2-C7	5.08	129.70	122.90
7	C	1308	NAG	C2-N2-C7	5.07	129.70	122.90
7	B	1303	NAG	C2-N2-C7	5.02	129.63	122.90
7	B	1304	NAG	C2-N2-C7	5.02	129.63	122.90
7	B	1309	NAG	C2-N2-C7	5.02	129.62	122.90
7	A	1307	NAG	C1-O5-C5	4.51	118.23	112.19
7	A	1301	NAG	C1-O5-C5	4.49	118.21	112.19
7	B	1308	NAG	C1-O5-C5	4.26	117.89	112.19
7	A	1309	NAG	C1-O5-C5	4.24	117.87	112.19
7	B	1302	NAG	C1-O5-C5	4.19	117.80	112.19
7	C	1307	NAG	C1-O5-C5	3.86	117.36	112.19
7	A	1303	NAG	C1-O5-C5	3.81	117.29	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	1303	NAG	C1-O5-C5	3.79	117.27	112.19
7	C	1304	NAG	C1-O5-C5	3.78	117.25	112.19
7	A	1308	NAG	C1-O5-C5	3.77	117.24	112.19
7	B	1306	NAG	C1-O5-C5	3.74	117.20	112.19
7	B	1303	NAG	C1-O5-C5	3.74	117.19	112.19
7	A	1304	NAG	C1-O5-C5	3.54	116.93	112.19
7	C	1308	NAG	C1-O5-C5	3.47	116.84	112.19
7	A	1306	NAG	C1-O5-C5	3.44	116.80	112.19
7	B	1310	NAG	C1-O5-C5	3.41	116.75	112.19
7	A	1302	NAG	C1-O5-C5	3.35	116.67	112.19
7	C	1309	NAG	C1-O5-C5	3.31	116.63	112.19
7	C	1301	NAG	C1-O5-C5	3.31	116.63	112.19
7	B	1308	NAG	C2-N2-C7	3.31	127.34	122.90
7	B	1305	NAG	C1-O5-C5	3.26	116.56	112.19
7	B	1307	NAG	C1-O5-C5	3.25	116.54	112.19
7	C	1306	NAG	C1-O5-C5	3.22	116.50	112.19
7	B	1304	NAG	C1-O5-C5	3.06	116.28	112.19
7	C	1305	NAG	C1-O5-C5	3.02	116.23	112.19
7	A	1307	NAG	C2-N2-C7	3.00	126.92	122.90
7	B	1309	NAG	C1-C2-N2	2.85	114.93	110.43
7	B	1303	NAG	C1-C2-N2	2.81	114.87	110.43
7	B	1301	NAG	C1-O5-C5	2.76	115.88	112.19
7	A	1308	NAG	C1-C2-N2	2.63	114.58	110.43
7	C	1308	NAG	C1-C2-N2	2.57	114.48	110.43
7	A	1305	NAG	C1-O5-C5	2.56	115.61	112.19
7	C	1302	NAG	C1-O5-C5	2.47	115.49	112.19
7	B	1304	NAG	C1-C2-N2	2.45	114.30	110.43
7	B	1309	NAG	C1-O5-C5	2.42	115.44	112.19

There are no chirality outliers.

All (65) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1306	NAG	O5-C5-C6-O6
7	C	1304	NAG	O5-C5-C6-O6
7	B	1302	NAG	O5-C5-C6-O6
7	B	1307	NAG	O5-C5-C6-O6
7	A	1308	NAG	O5-C5-C6-O6
7	A	1307	NAG	O5-C5-C6-O6
7	B	1308	NAG	O5-C5-C6-O6
7	C	1303	NAG	O5-C5-C6-O6
7	B	1307	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	C	1304	NAG	C4-C5-C6-O6
7	B	1310	NAG	O5-C5-C6-O6
7	C	1307	NAG	O5-C5-C6-O6
7	A	1308	NAG	C4-C5-C6-O6
7	A	1309	NAG	O5-C5-C6-O6
7	B	1303	NAG	O5-C5-C6-O6
7	A	1305	NAG	O5-C5-C6-O6
7	C	1308	NAG	C4-C5-C6-O6
7	C	1308	NAG	O5-C5-C6-O6
7	B	1301	NAG	O5-C5-C6-O6
7	C	1306	NAG	O5-C5-C6-O6
7	B	1309	NAG	O5-C5-C6-O6
7	C	1301	NAG	O5-C5-C6-O6
7	A	1306	NAG	C4-C5-C6-O6
7	A	1303	NAG	O5-C5-C6-O6
7	A	1302	NAG	O5-C5-C6-O6
7	A	1305	NAG	C4-C5-C6-O6
7	B	1302	NAG	C4-C5-C6-O6
7	A	1302	NAG	C4-C5-C6-O6
7	C	1303	NAG	C4-C5-C6-O6
7	C	1307	NAG	C4-C5-C6-O6
7	A	1303	NAG	C4-C5-C6-O6
7	A	1307	NAG	C4-C5-C6-O6
7	B	1303	NAG	C4-C5-C6-O6
7	B	1308	NAG	C4-C5-C6-O6
7	C	1306	NAG	C4-C5-C6-O6
7	C	1309	NAG	O5-C5-C6-O6
7	B	1309	NAG	C4-C5-C6-O6
7	A	1307	NAG	C8-C7-N2-C2
7	A	1307	NAG	O7-C7-N2-C2
7	B	1308	NAG	C8-C7-N2-C2
7	B	1308	NAG	O7-C7-N2-C2
7	A	1301	NAG	O5-C5-C6-O6
7	A	1301	NAG	C4-C5-C6-O6
7	B	1310	NAG	C4-C5-C6-O6
7	B	1306	NAG	O5-C5-C6-O6
7	B	1306	NAG	C4-C5-C6-O6
7	B	1304	NAG	O5-C5-C6-O6
7	B	1301	NAG	C4-C5-C6-O6
7	C	1302	NAG	C4-C5-C6-O6
7	A	1309	NAG	C4-C5-C6-O6
7	C	1305	NAG	C4-C5-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
7	C	1305	NAG	O5-C5-C6-O6
7	C	1302	NAG	O5-C5-C6-O6
7	C	1301	NAG	C4-C5-C6-O6
7	B	1303	NAG	C3-C2-N2-C7
7	B	1304	NAG	C3-C2-N2-C7
7	B	1309	NAG	C3-C2-N2-C7
7	C	1309	NAG	C4-C5-C6-O6
7	A	1308	NAG	C1-C2-N2-C7
7	B	1303	NAG	C1-C2-N2-C7
7	B	1304	NAG	C1-C2-N2-C7
7	B	1309	NAG	C1-C2-N2-C7
7	C	1308	NAG	C1-C2-N2-C7
7	A	1308	NAG	C3-C2-N2-C7
7	C	1308	NAG	C3-C2-N2-C7

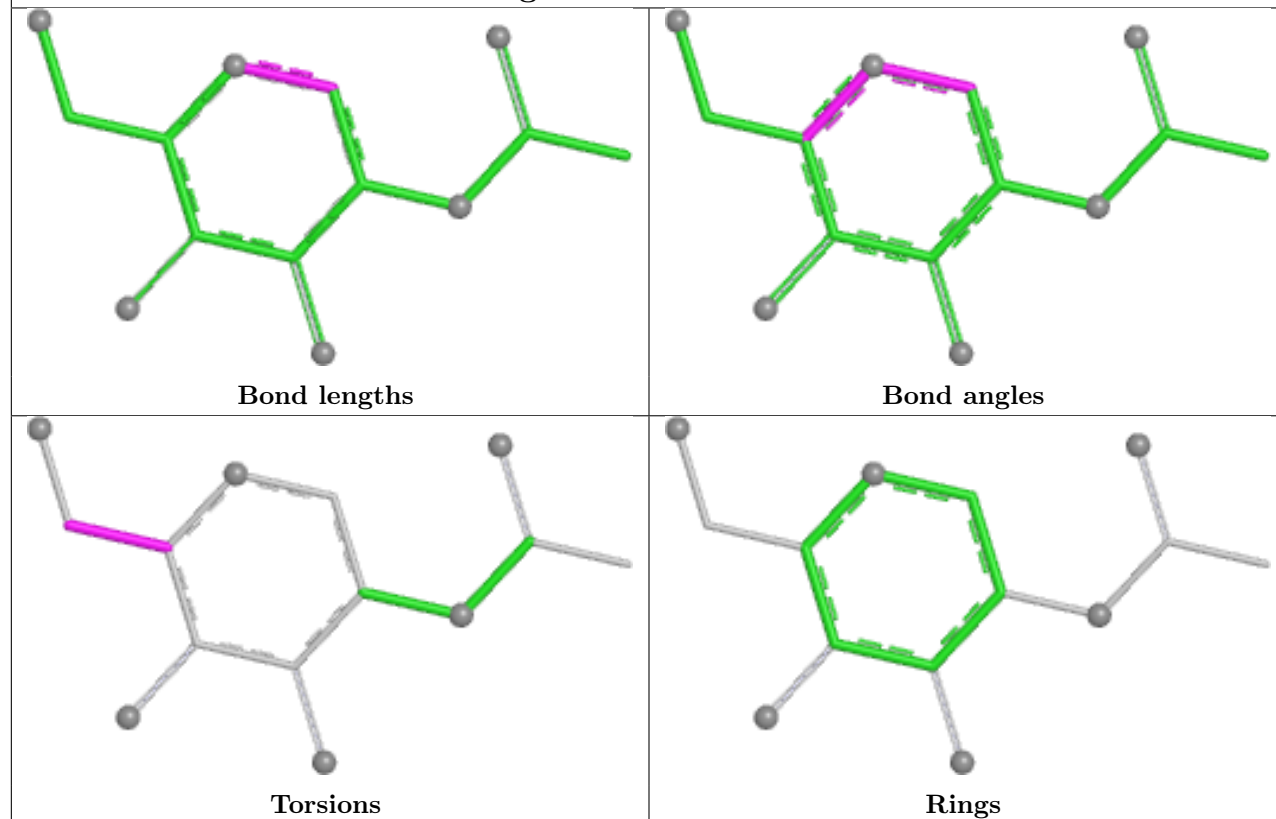
There are no ring outliers.

3 monomers are involved in 3 short contacts:

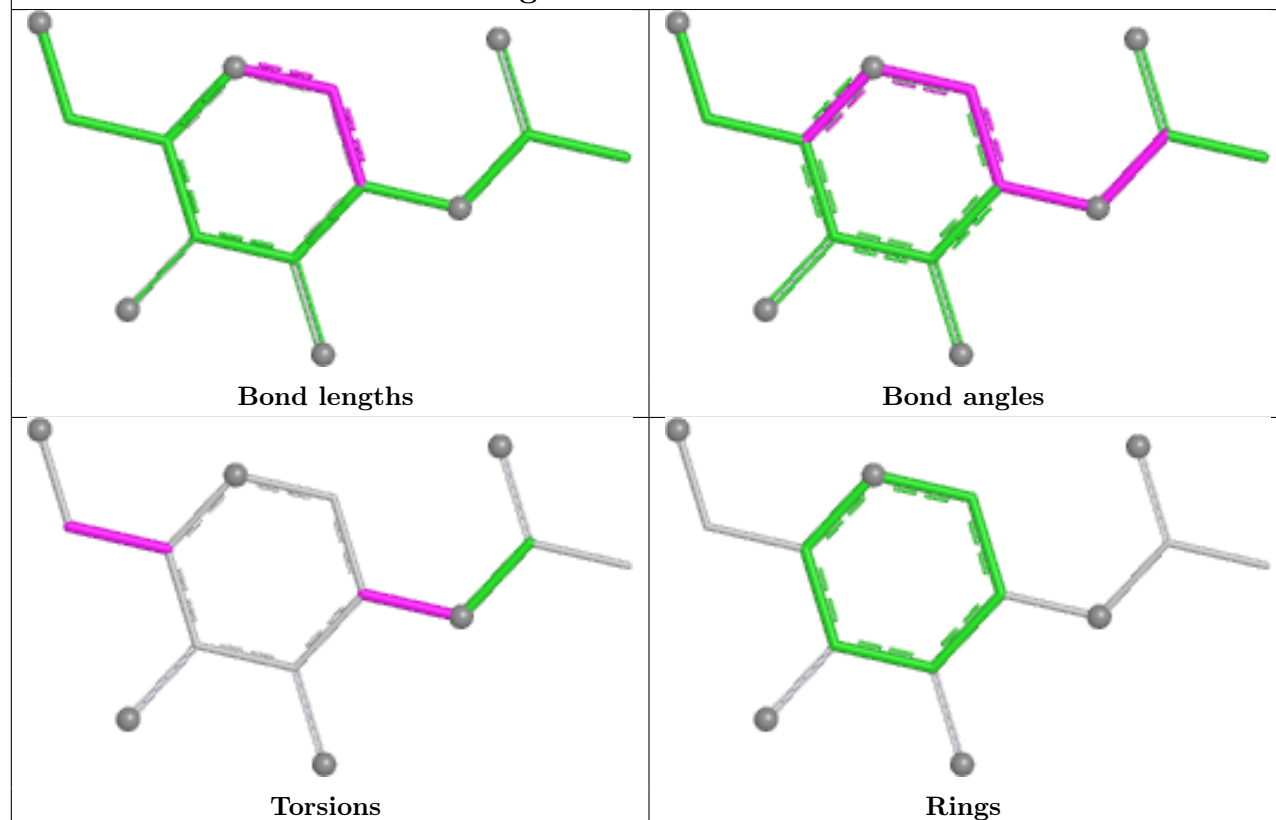
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	1306	NAG	1	0
7	B	1305	NAG	1	0
7	B	1308	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

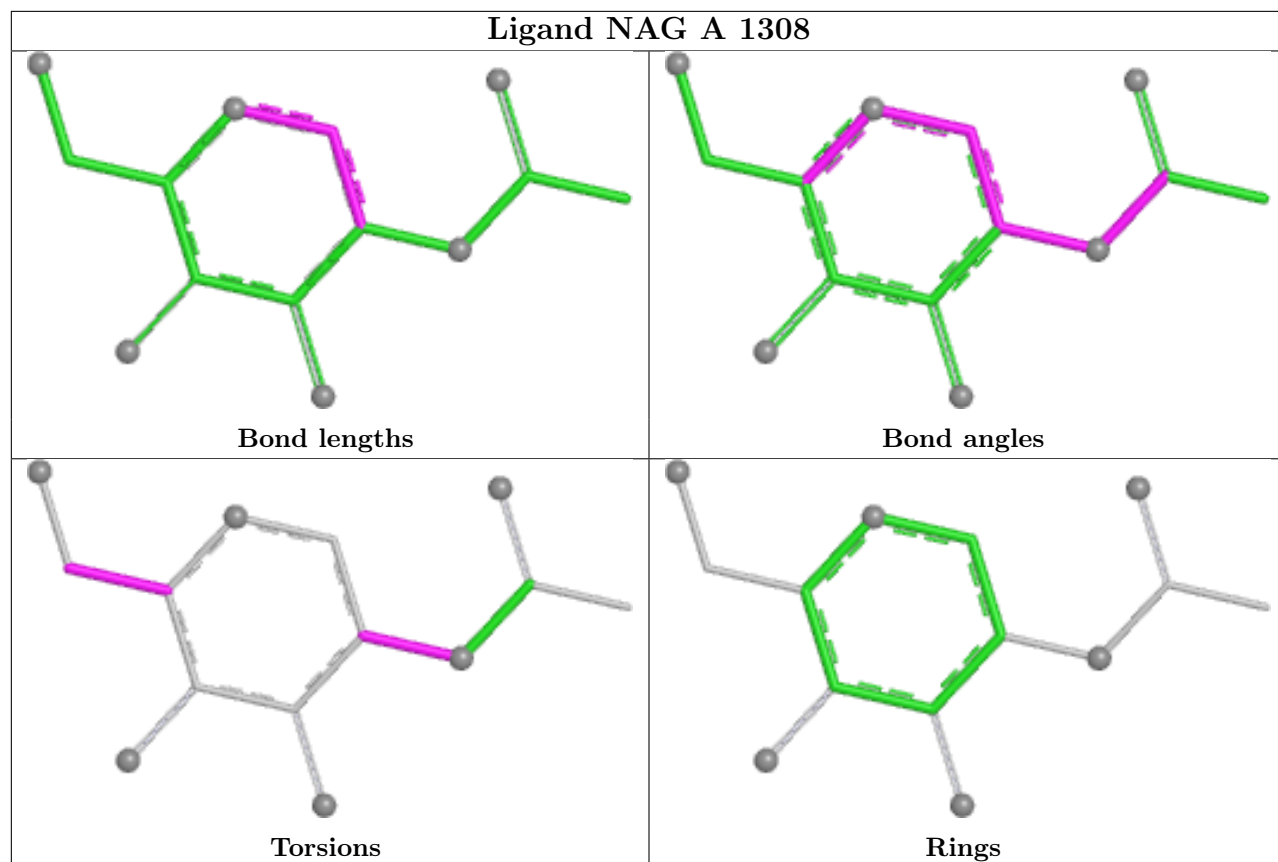
Ligand NAG C 1306



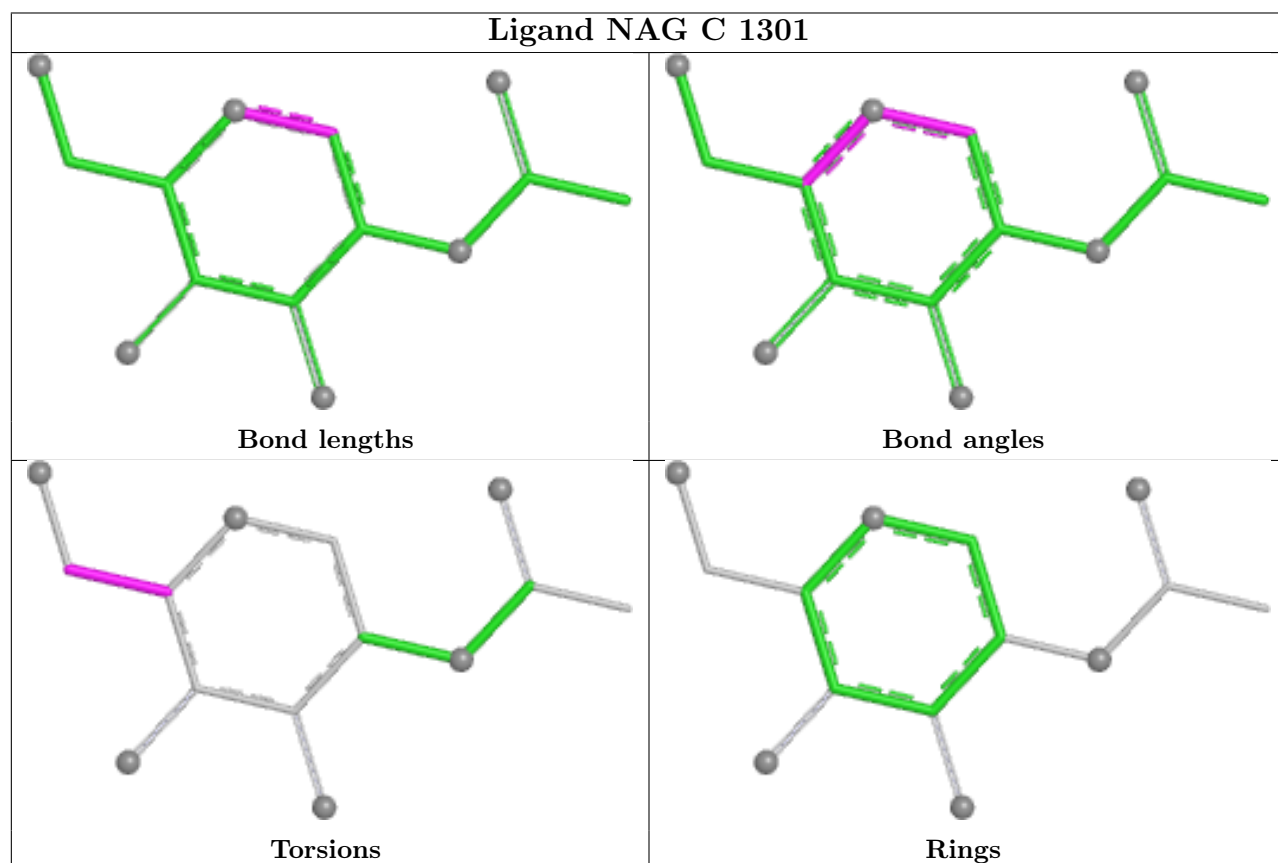
Ligand NAG B 1304

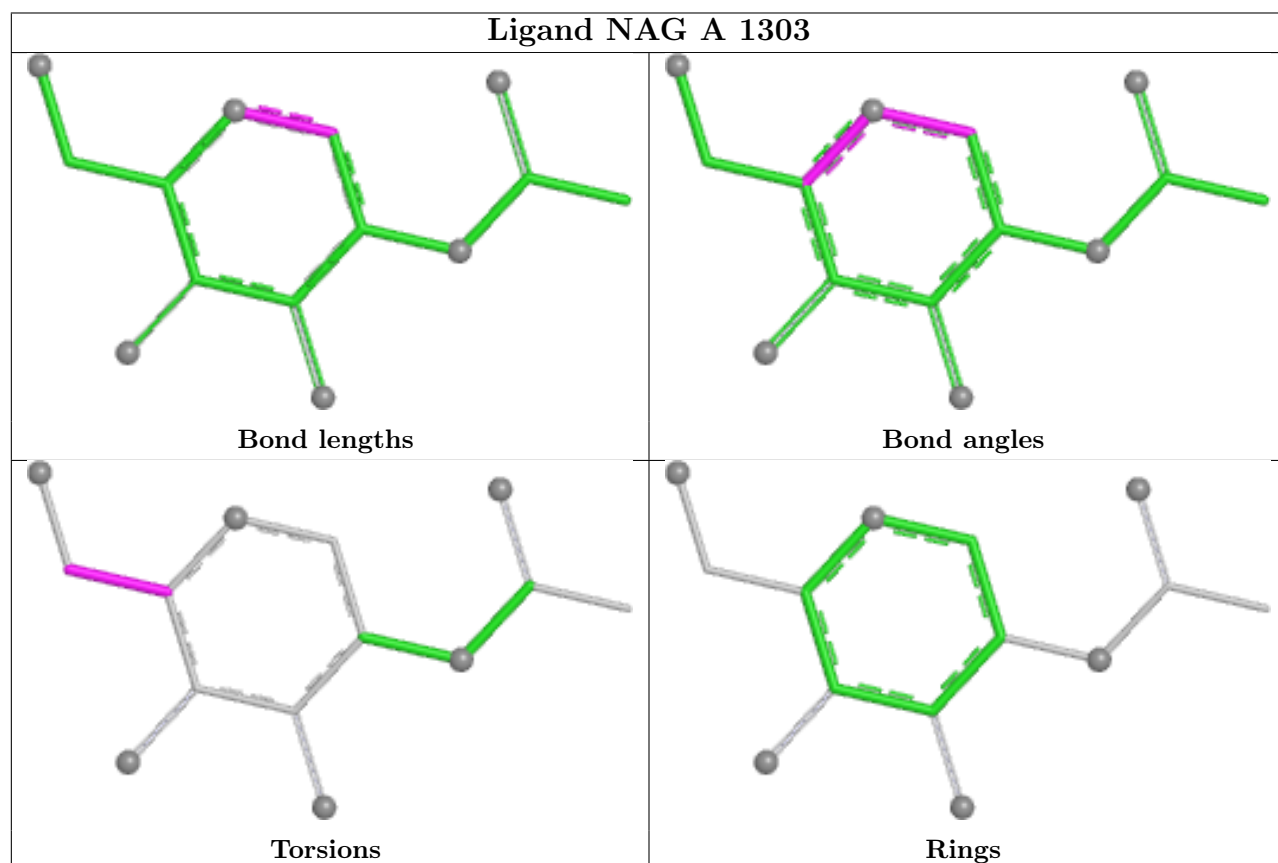
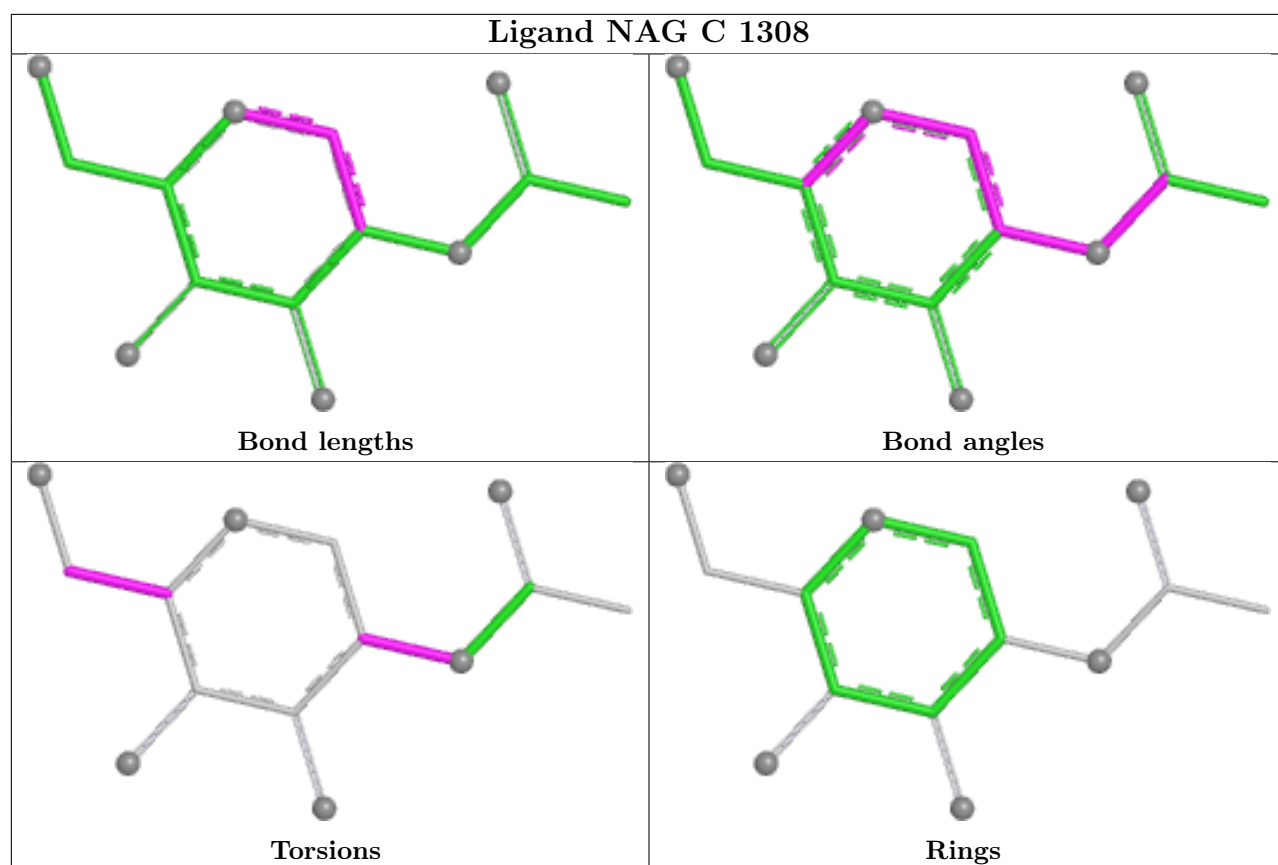


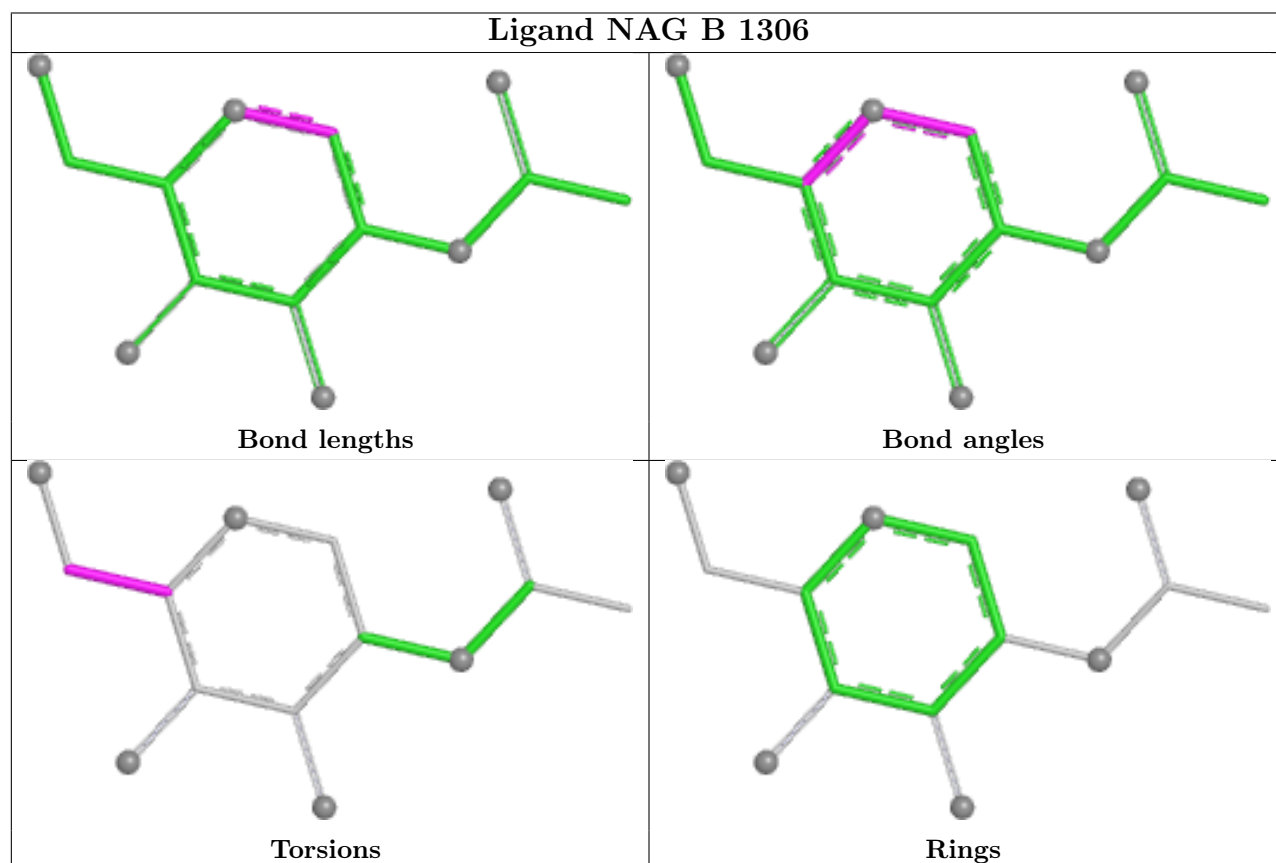
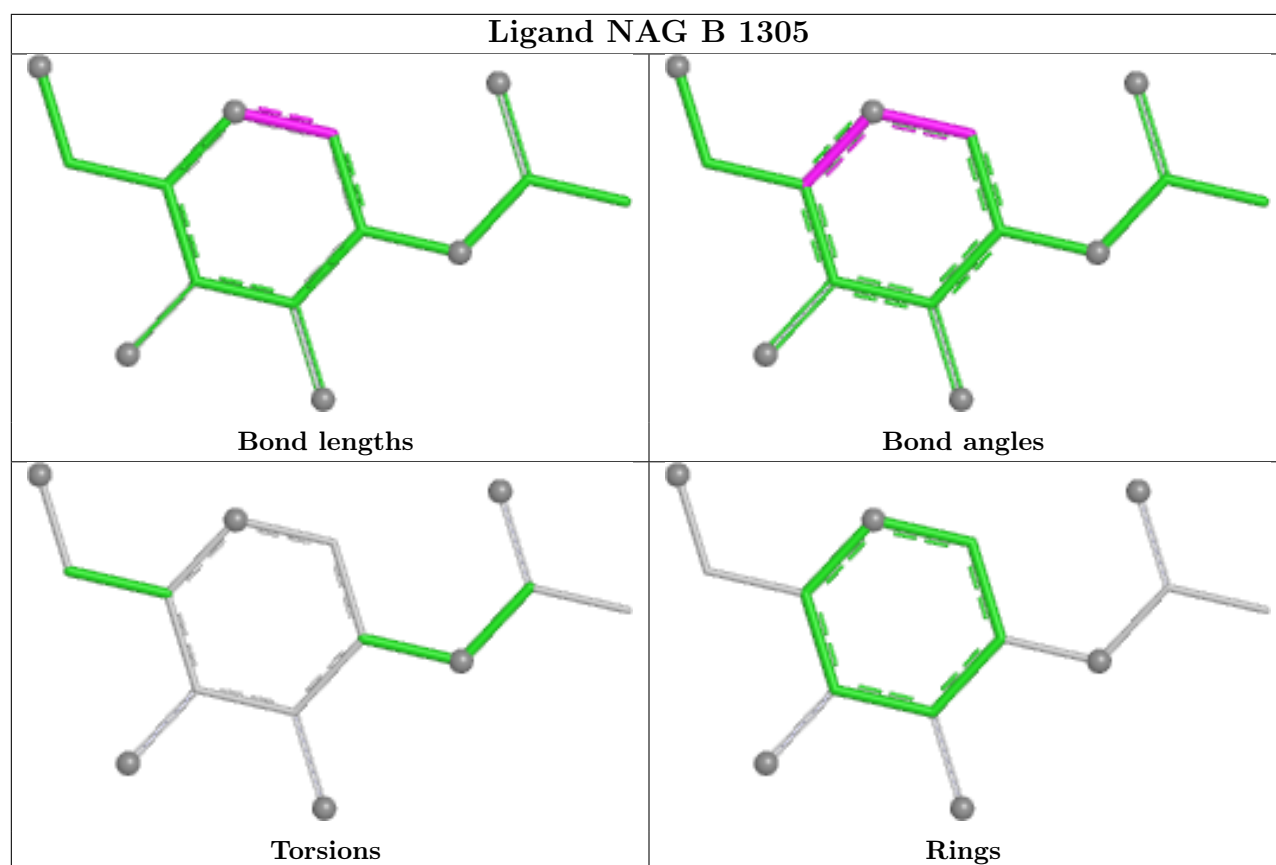
Ligand NAG A 1308

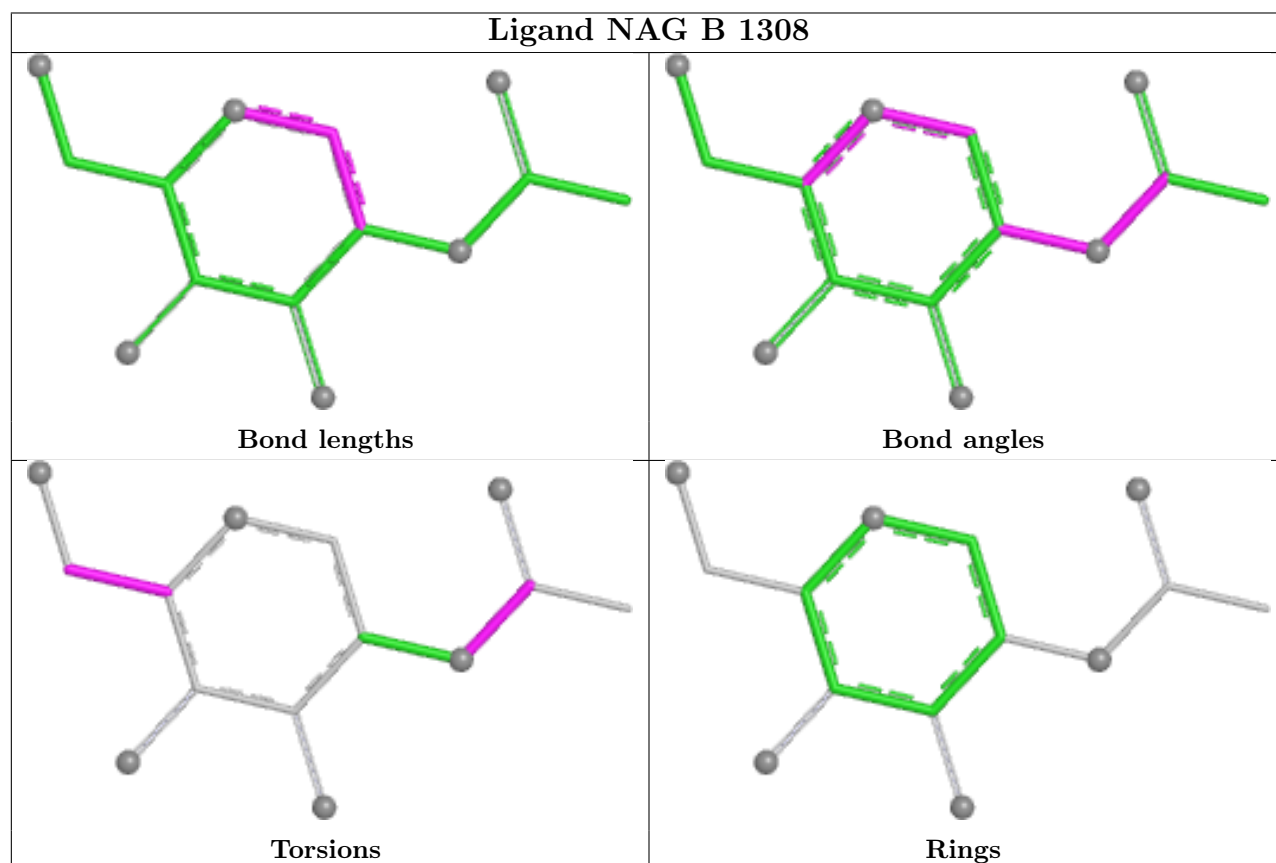
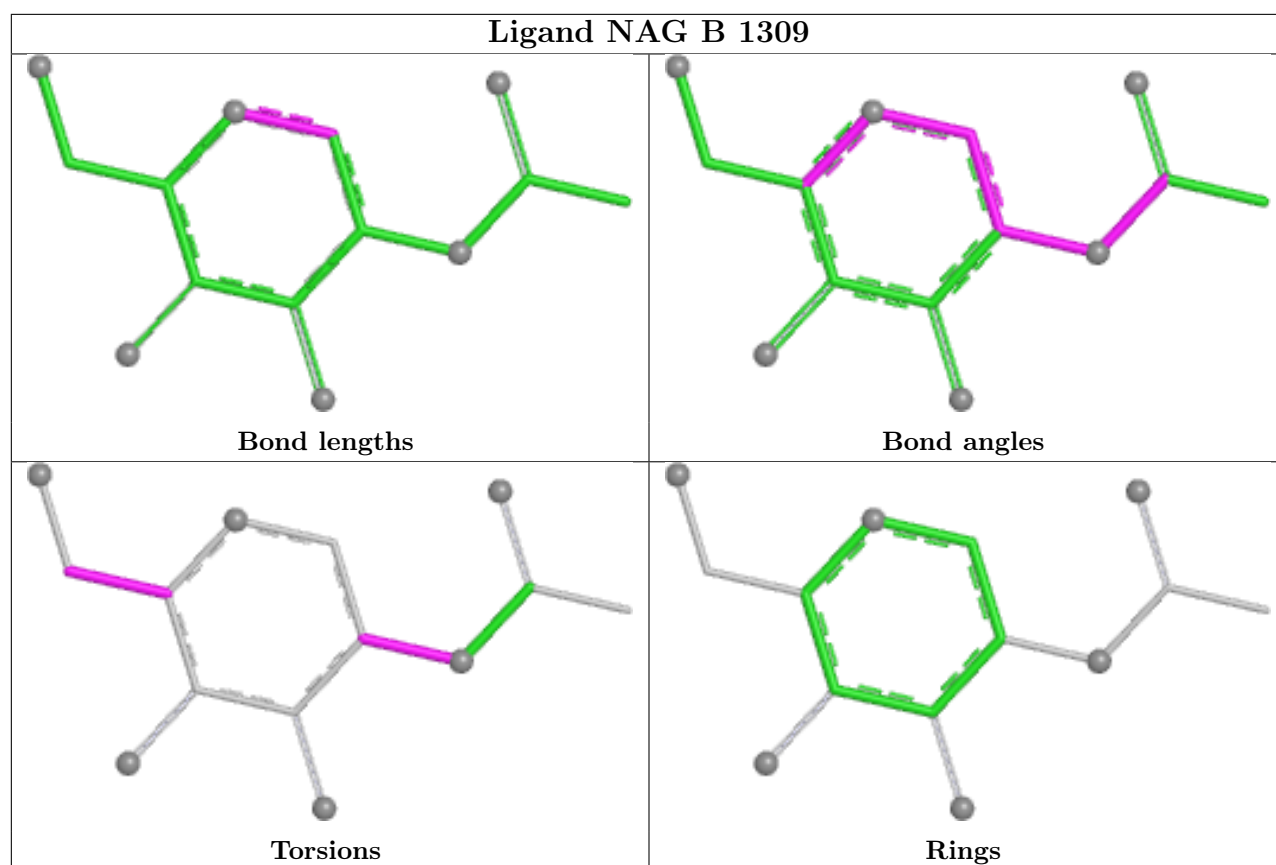


Ligand NAG C 1301

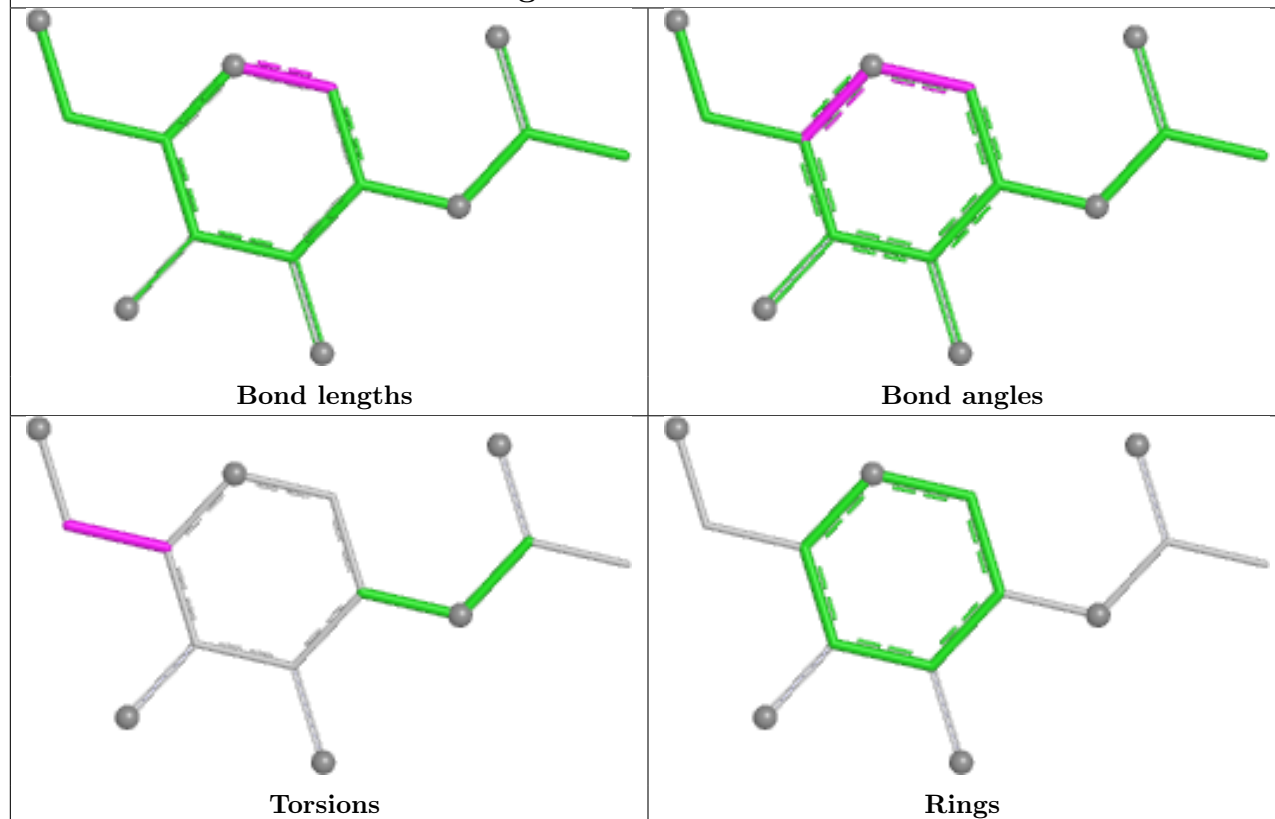




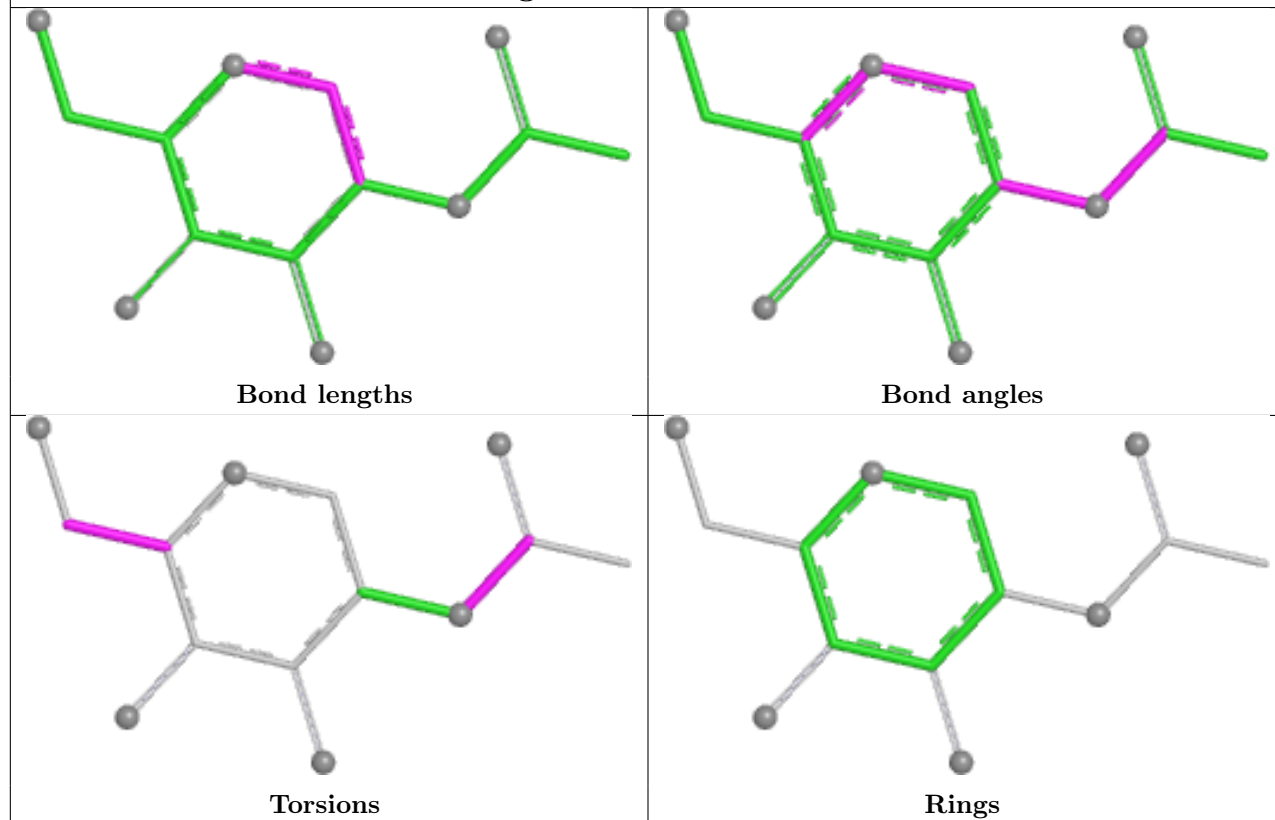




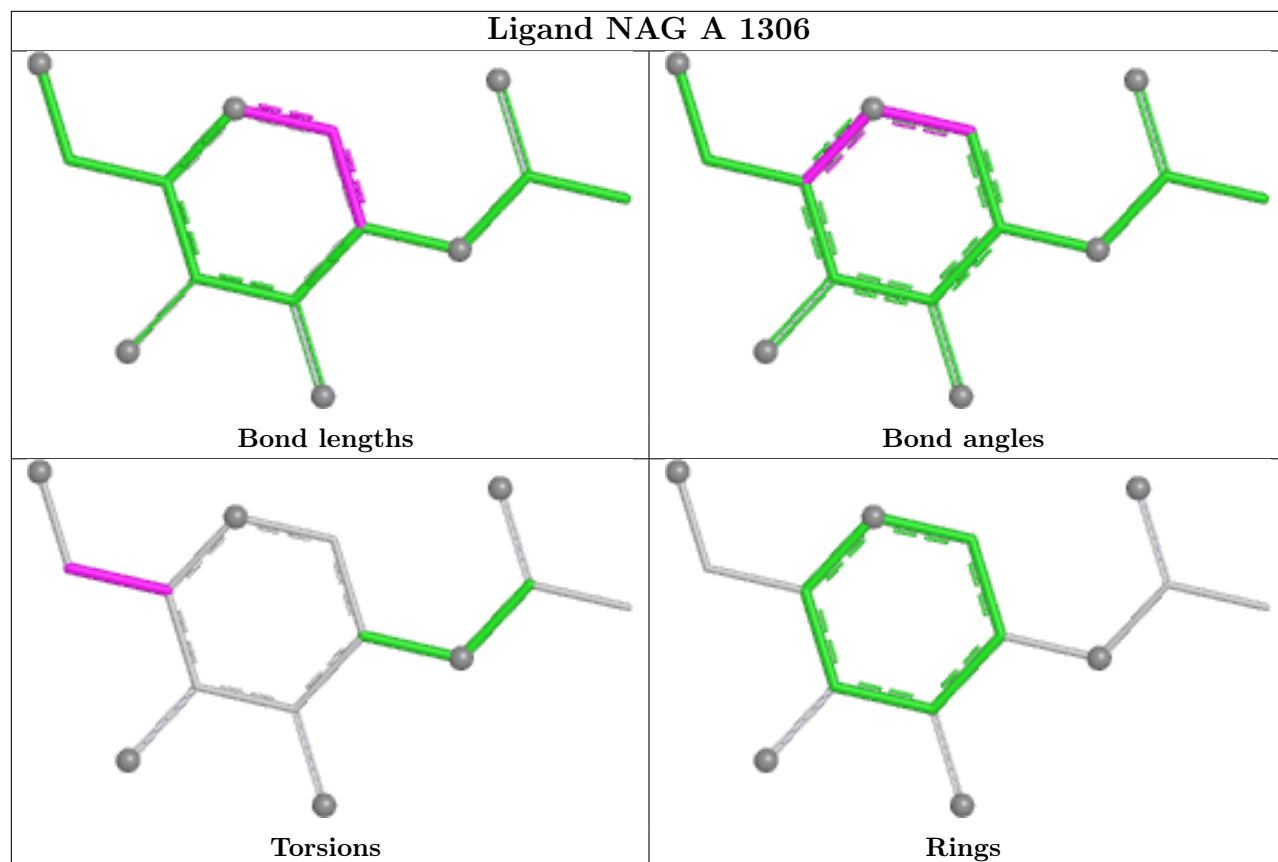
Ligand NAG C 1305



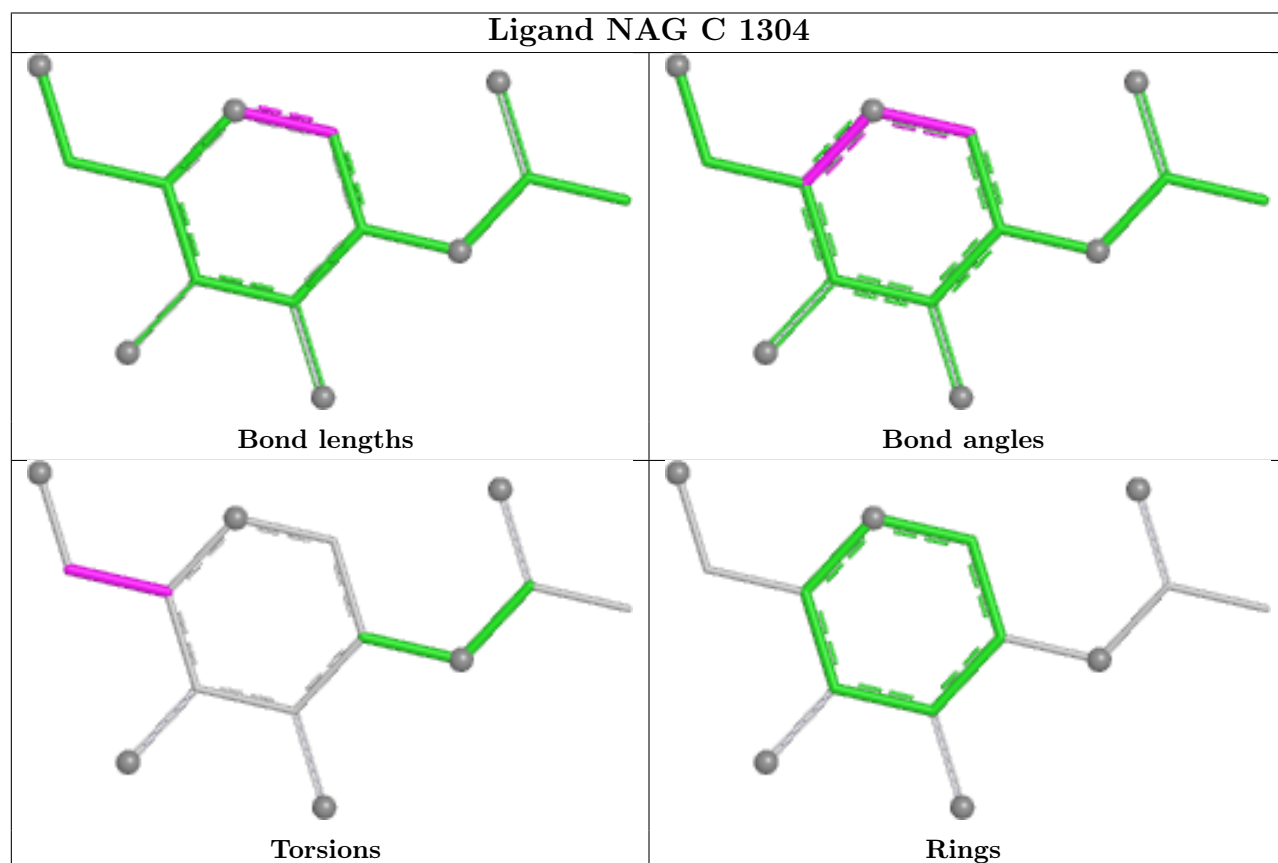
Ligand NAG A 1307



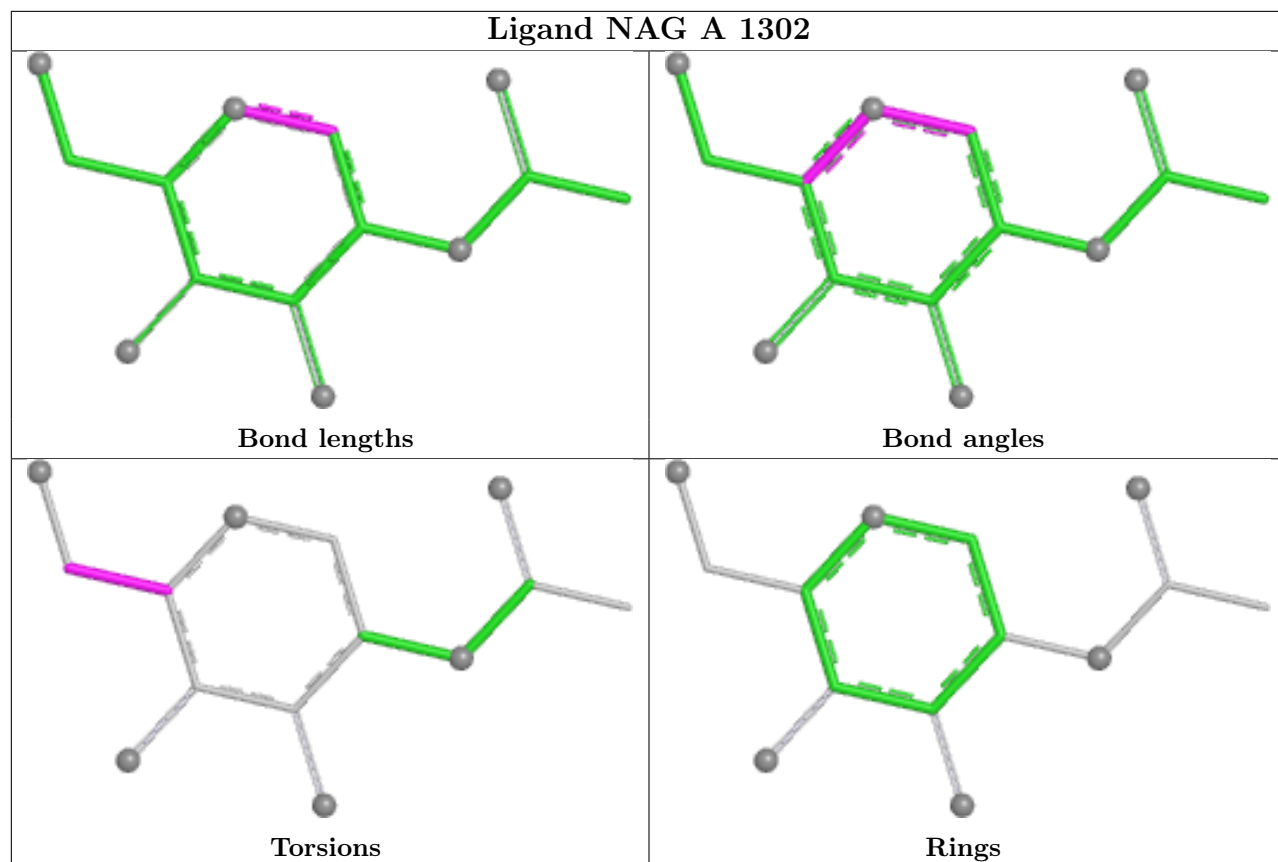
Ligand NAG A 1306



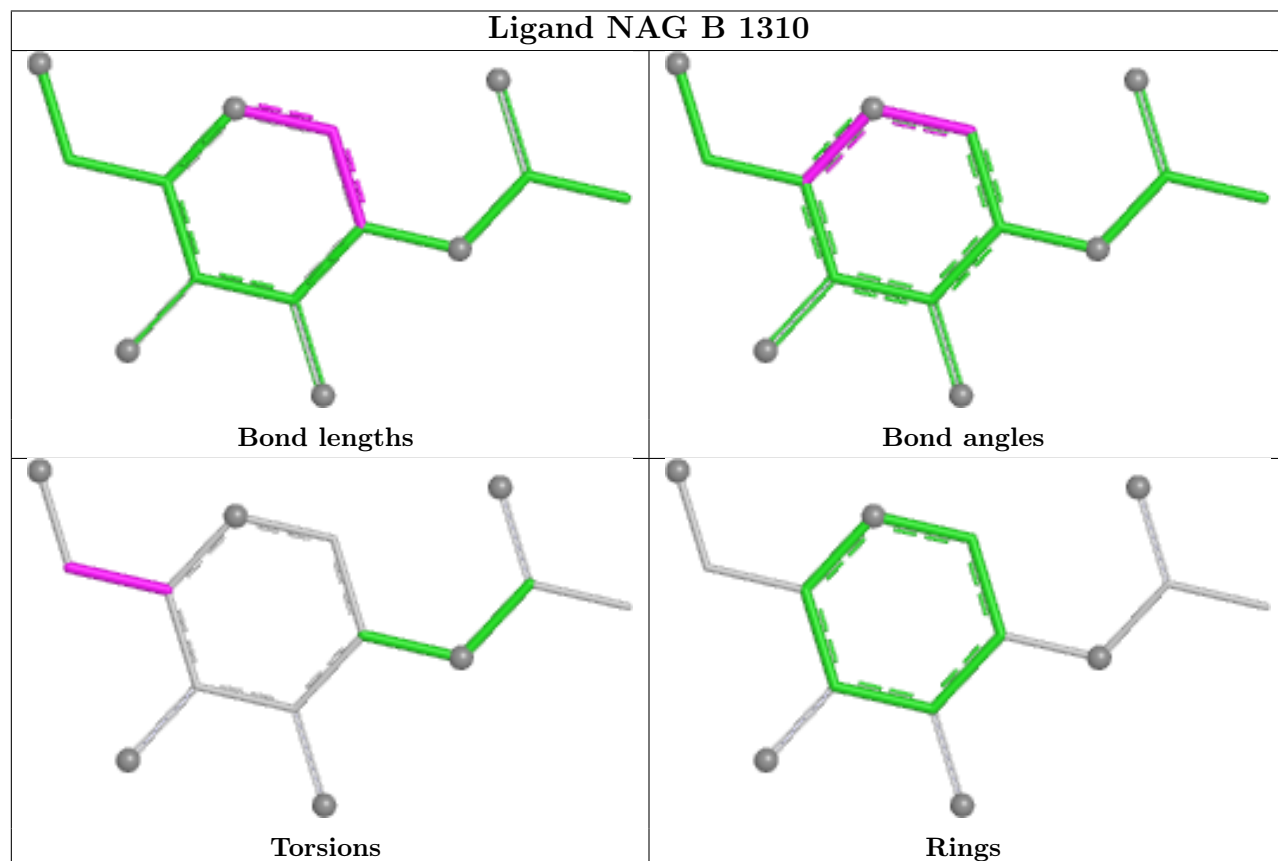
Ligand NAG C 1304



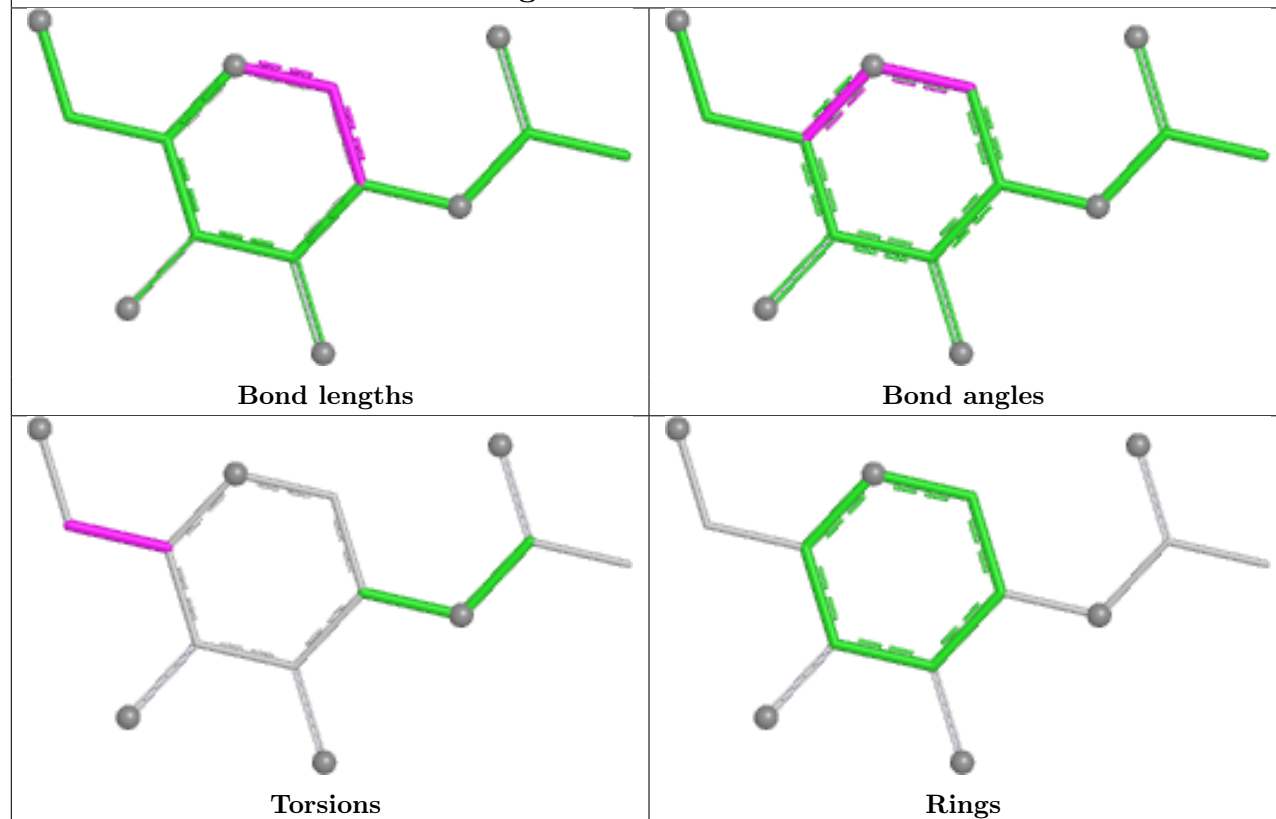
Ligand NAG A 1302



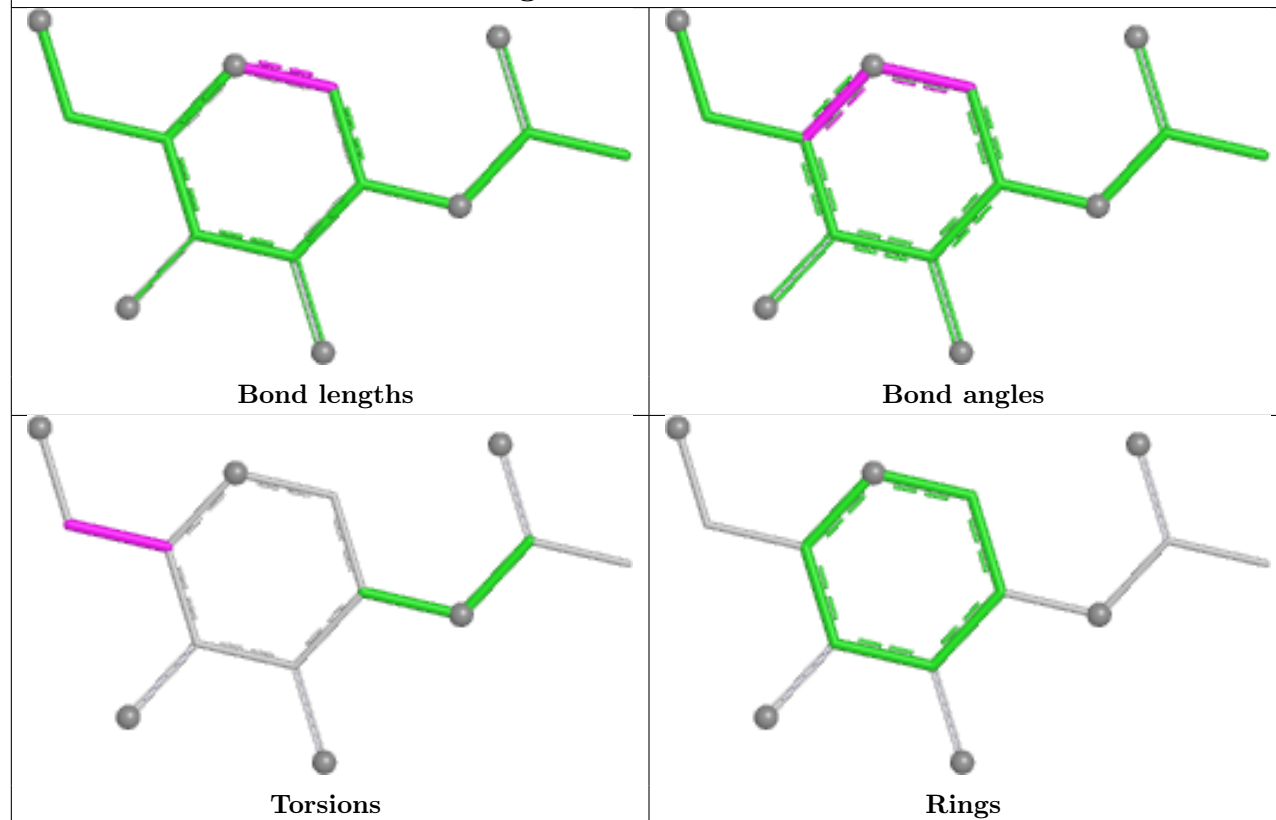
Ligand NAG B 1310

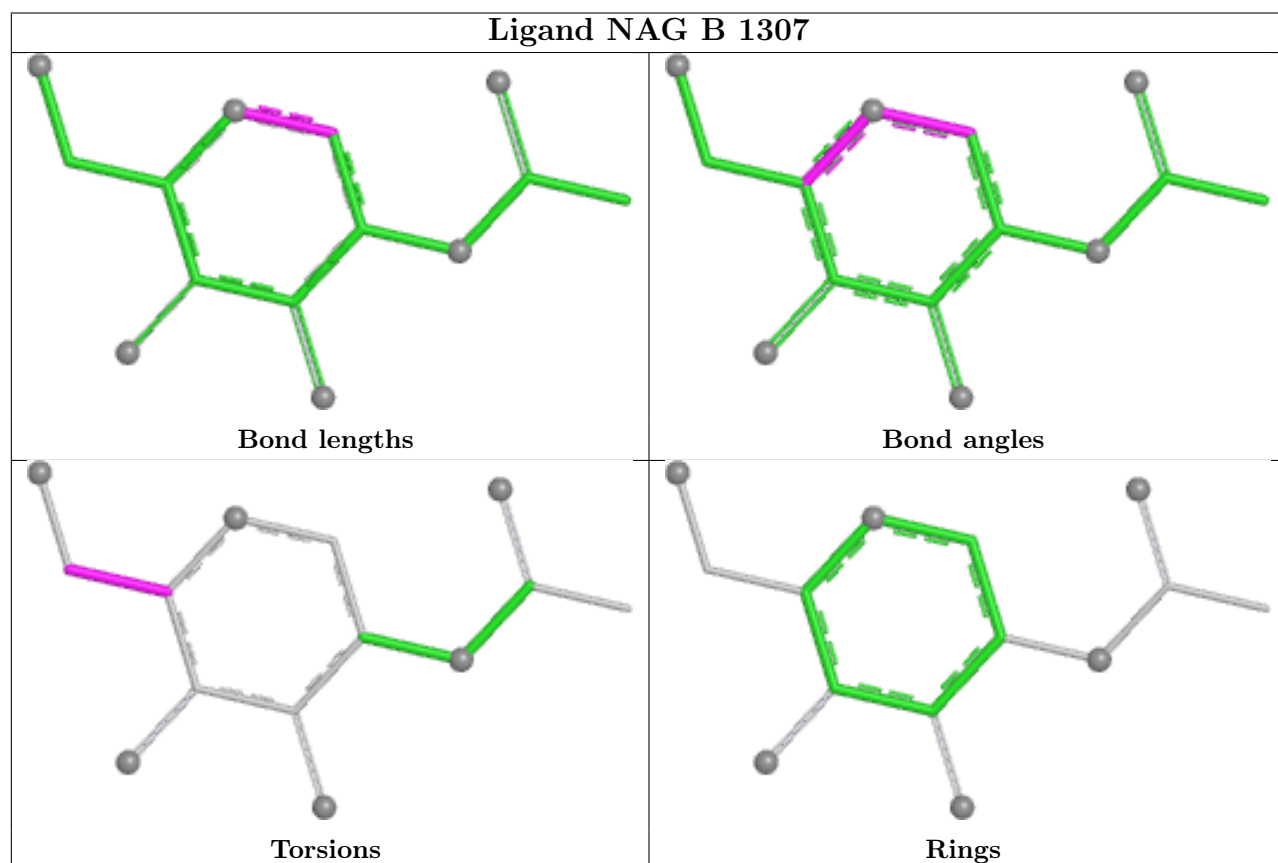
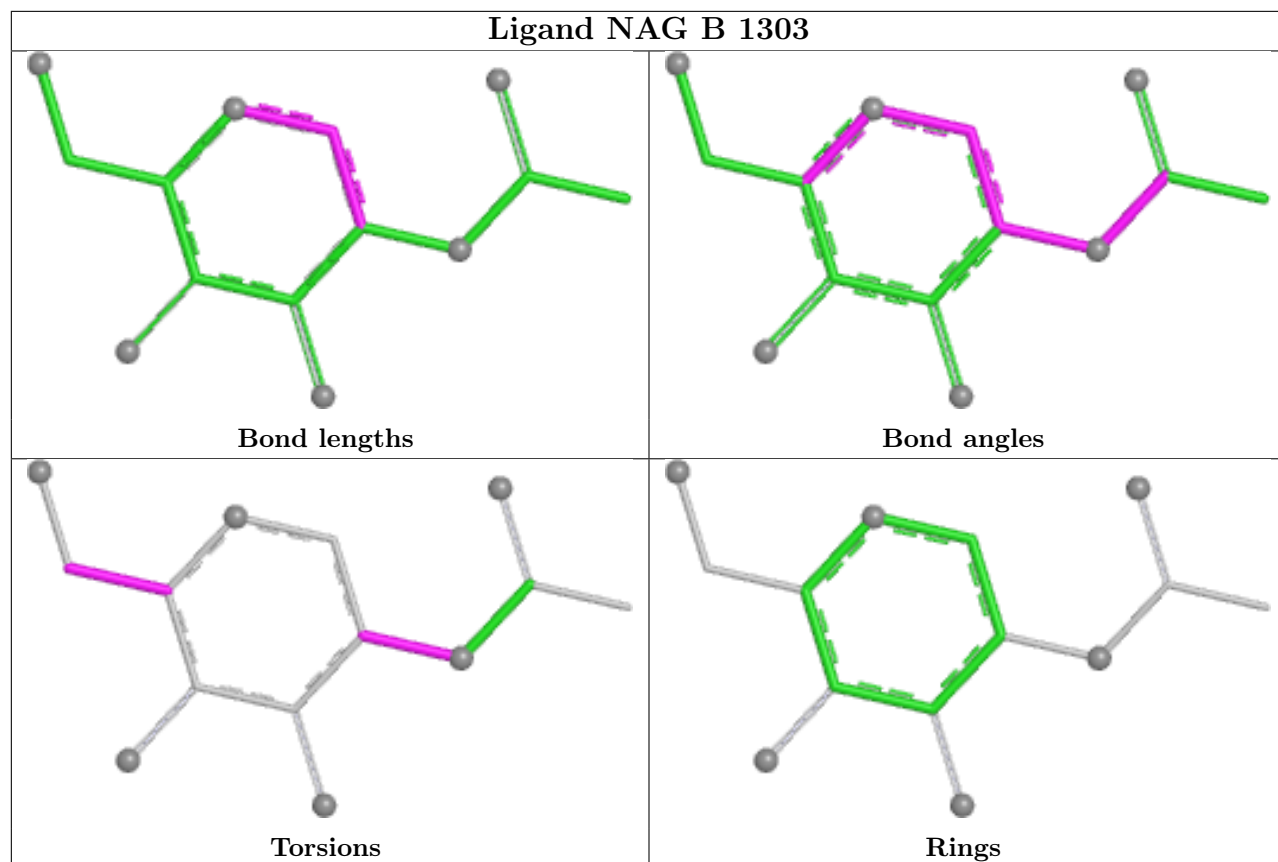


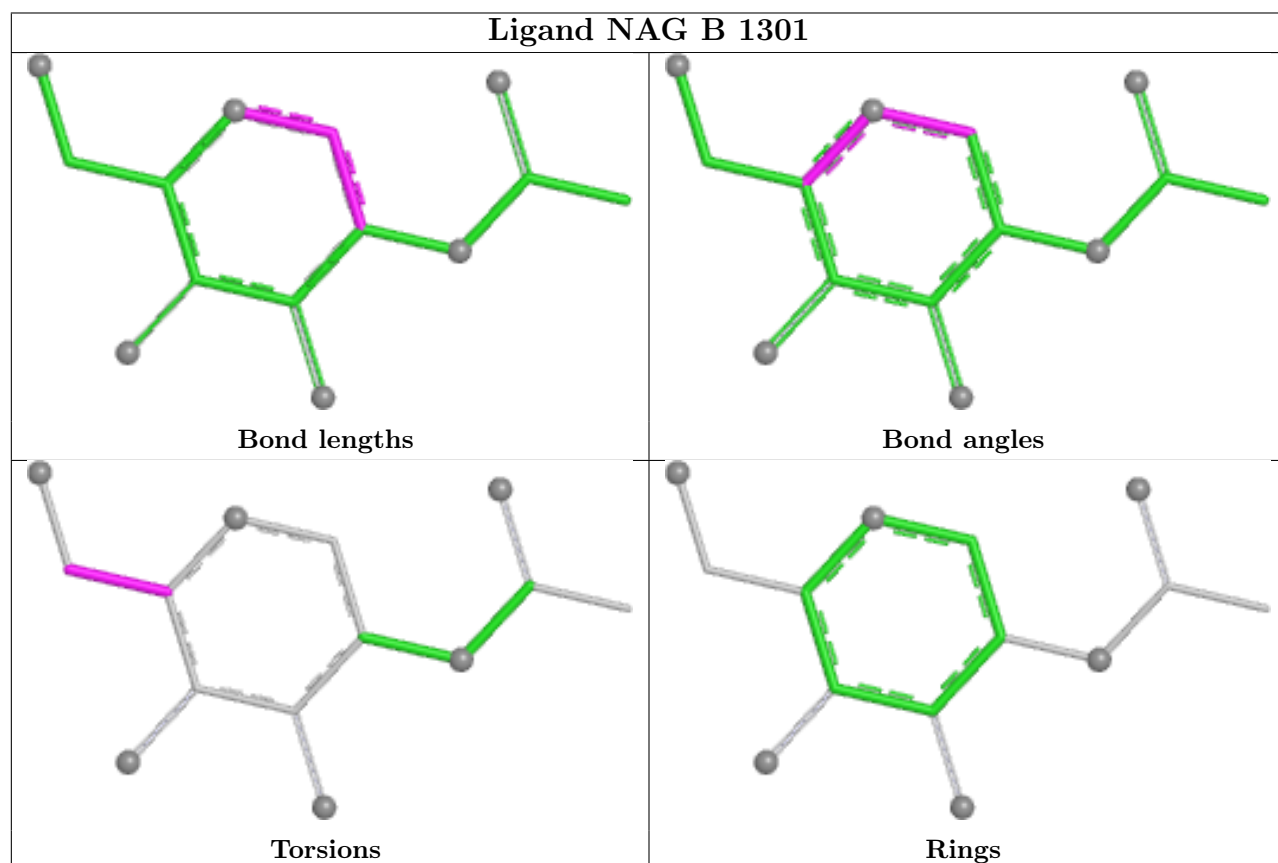
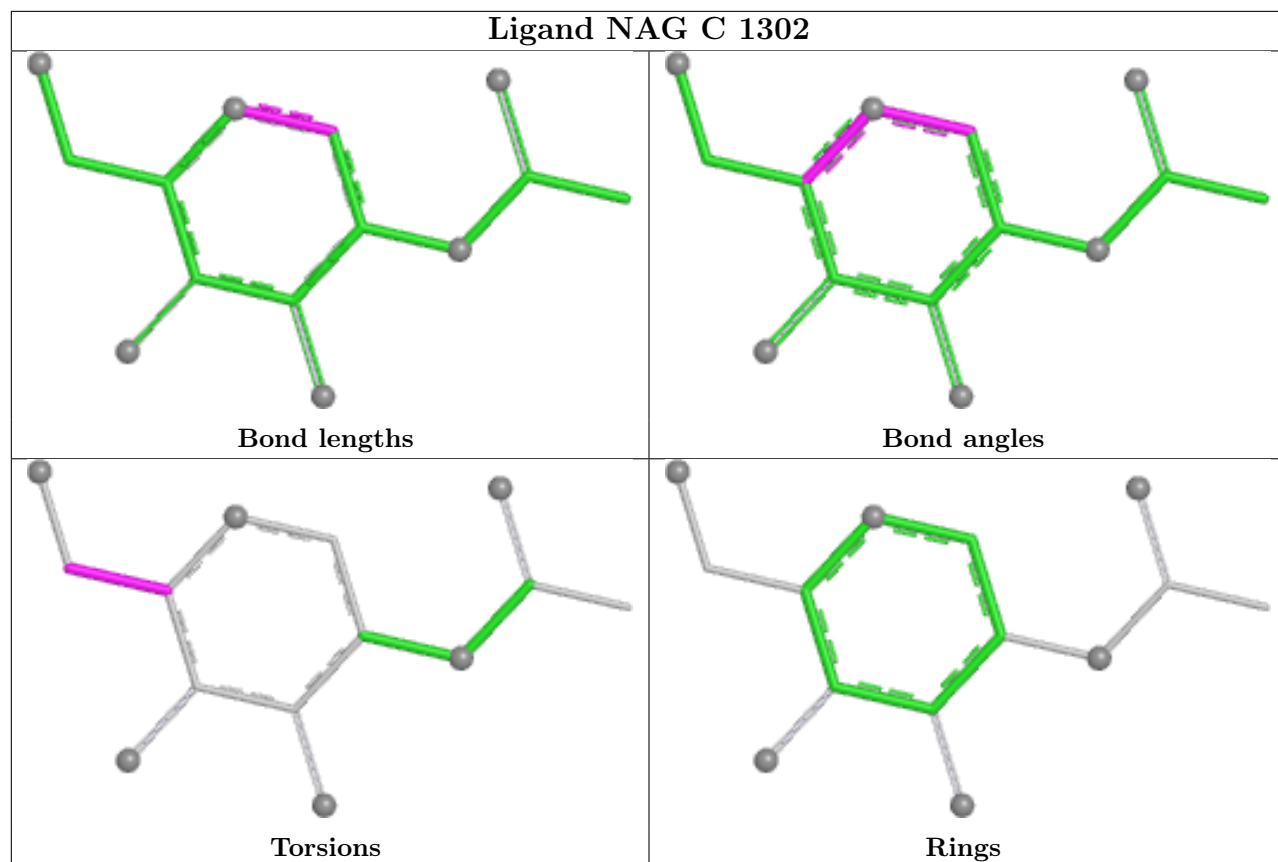
Ligand NAG A 1309



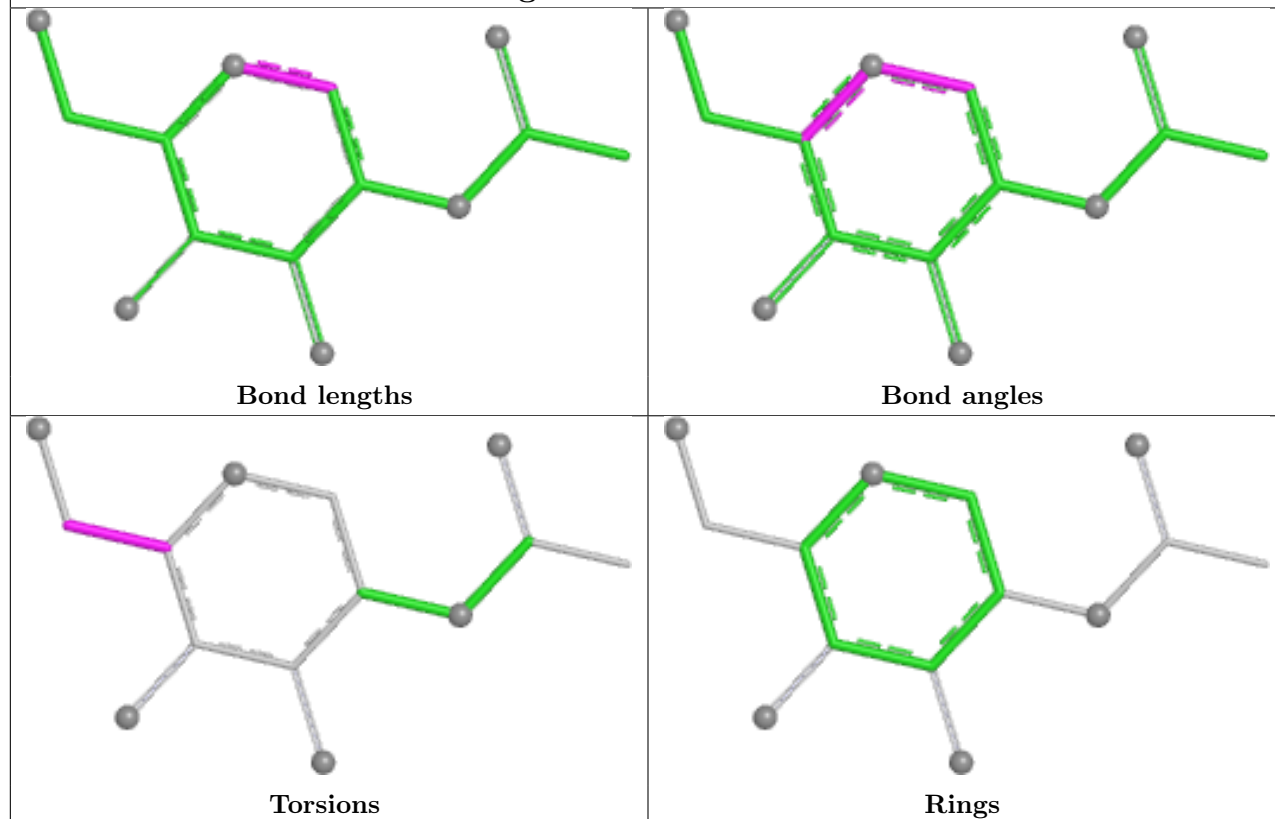
Ligand NAG A 1301



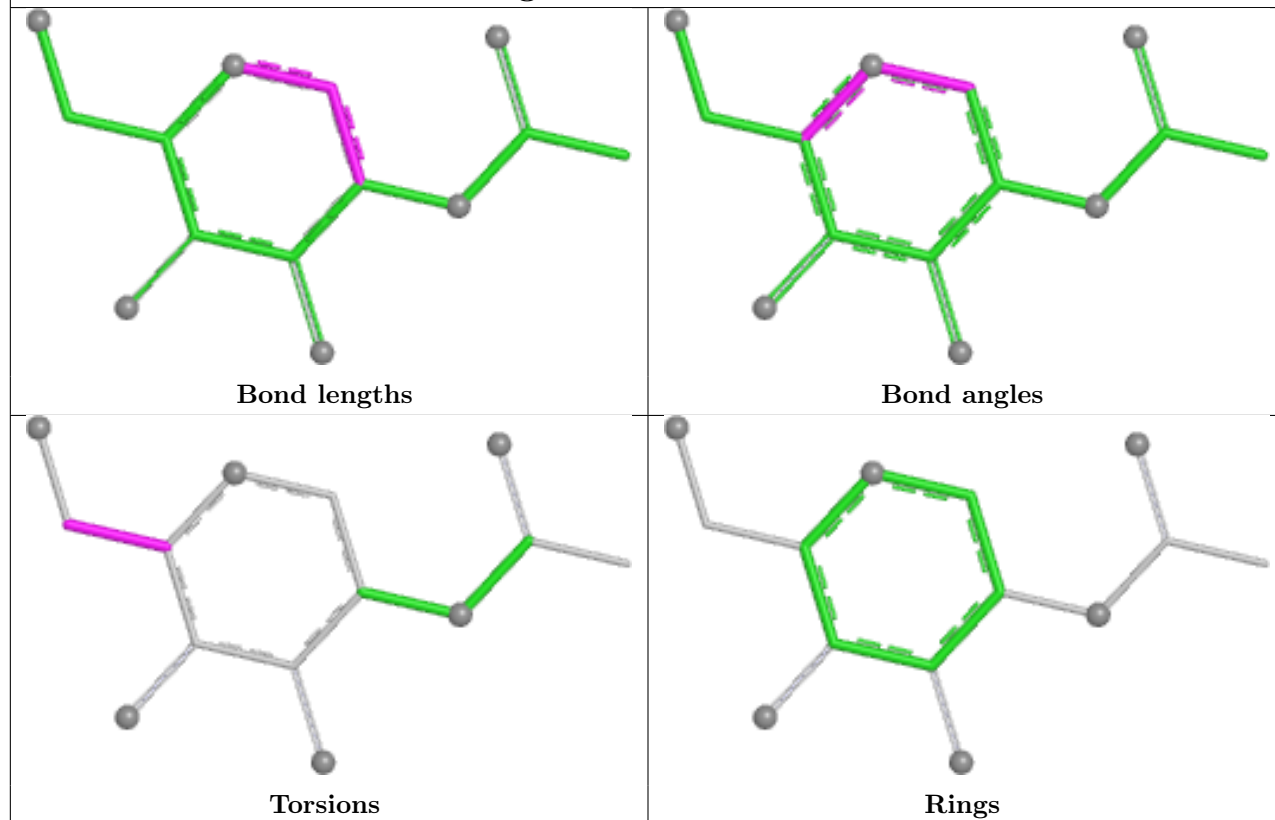


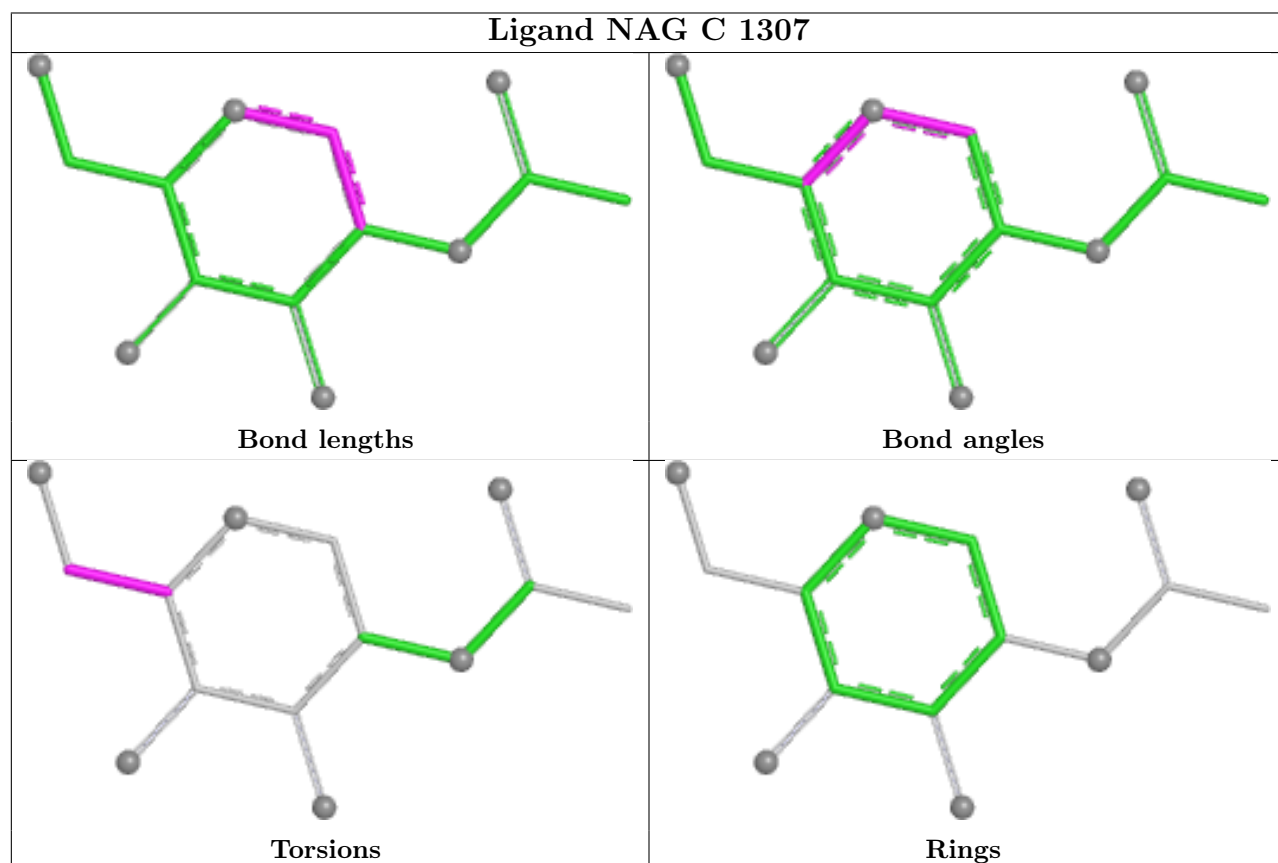
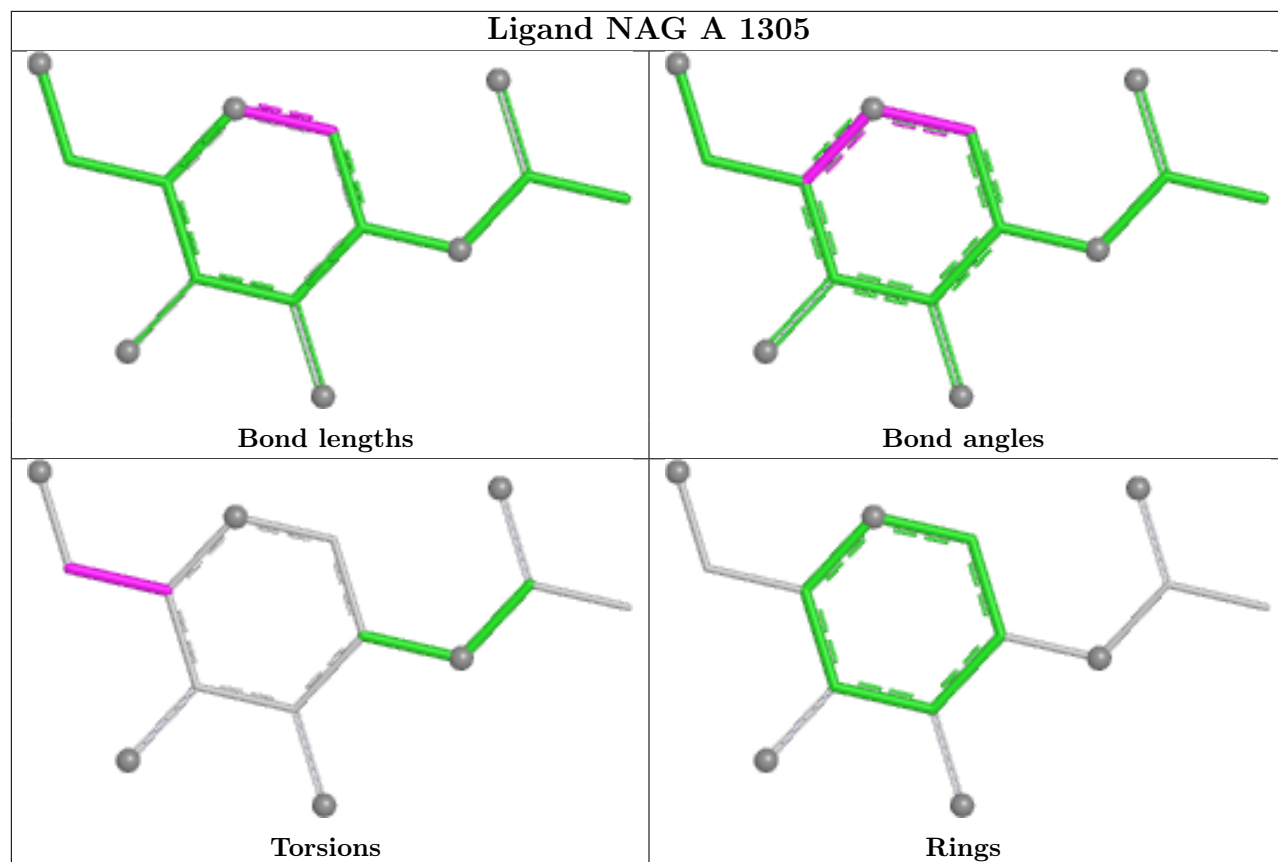


Ligand NAG B 1302

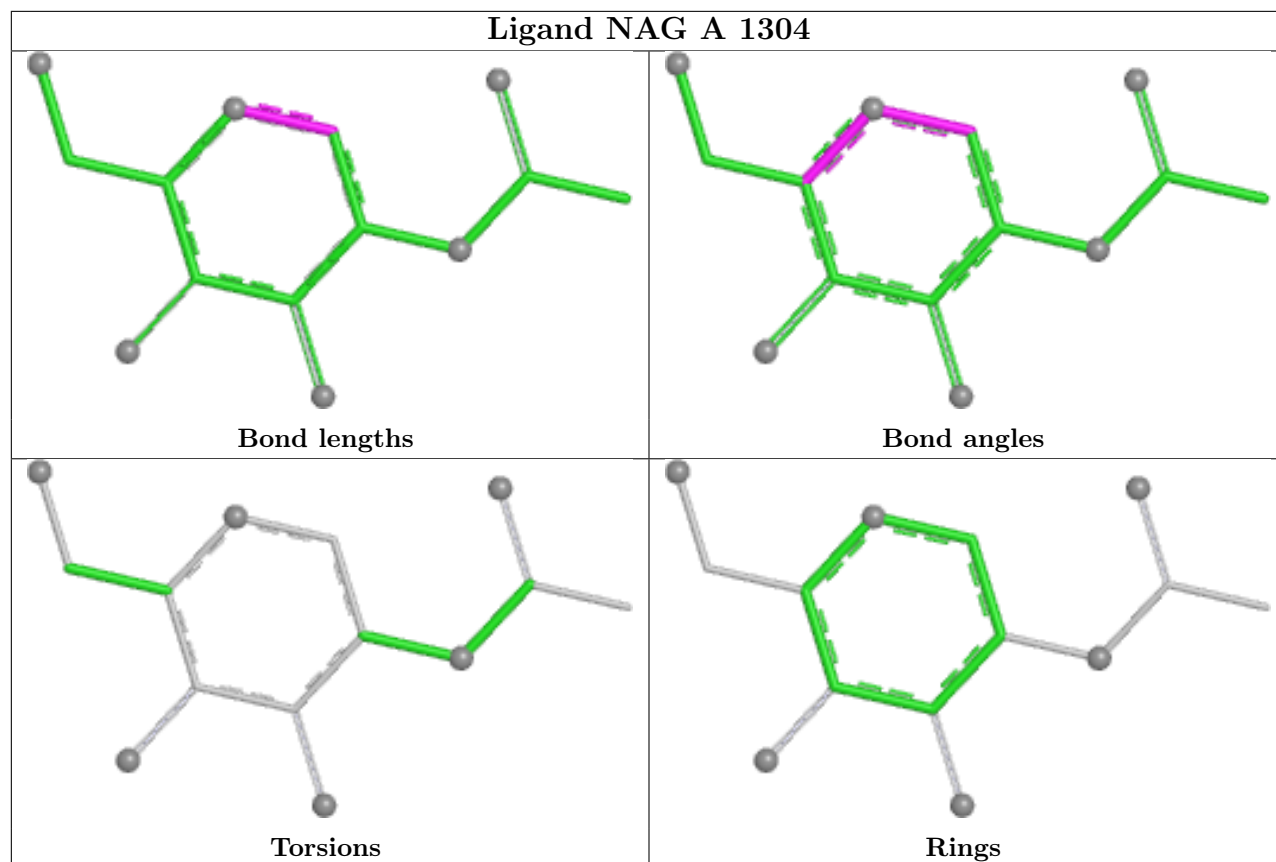


Ligand NAG C 1303

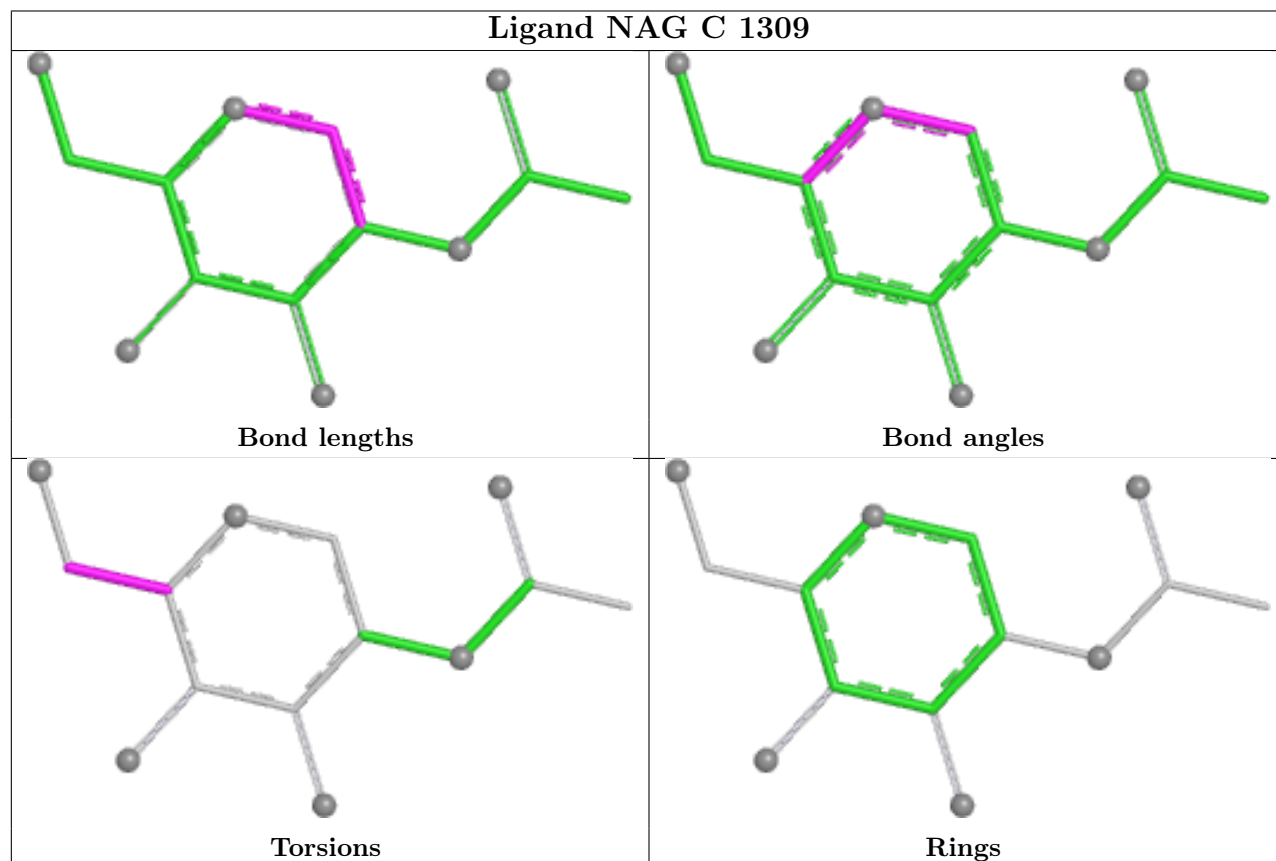




Ligand NAG A 1304



Ligand NAG C 1309



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

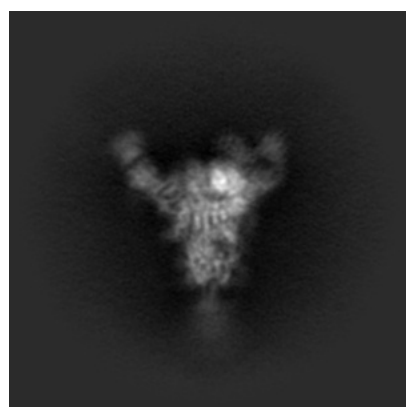
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24075. These allow visual inspection of the internal detail of the map and identification of artifacts.

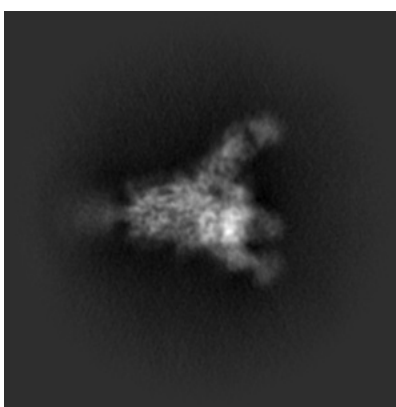
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

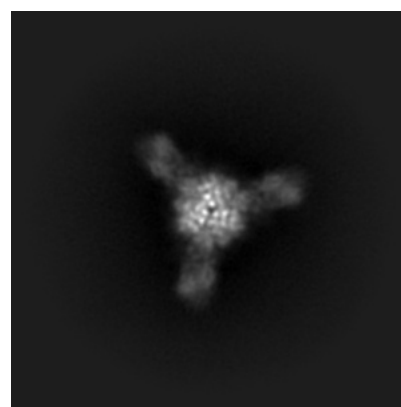
6.1.1 Primary map



X



Y

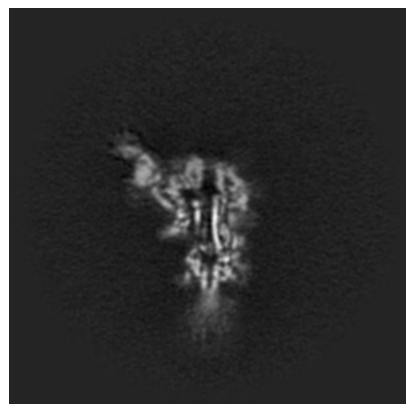


Z

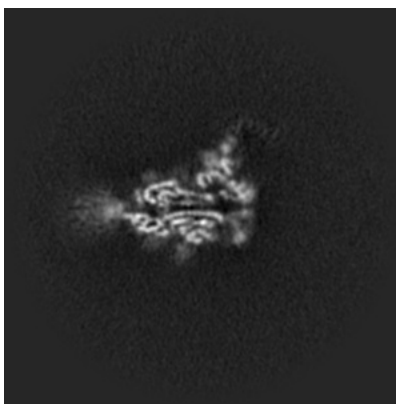
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

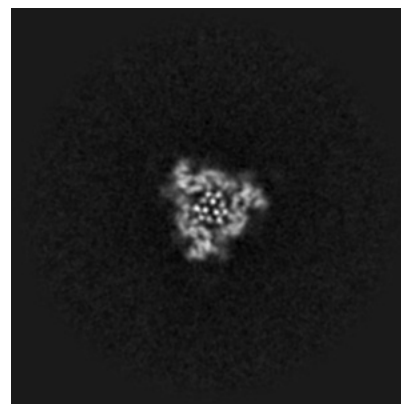
6.2.1 Primary map



X Index: 224



Y Index: 224

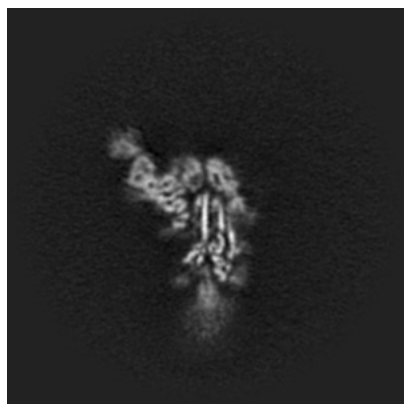


Z Index: 224

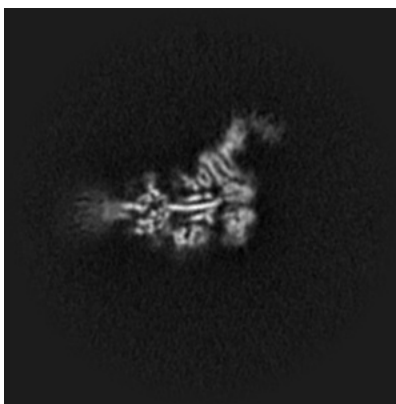
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

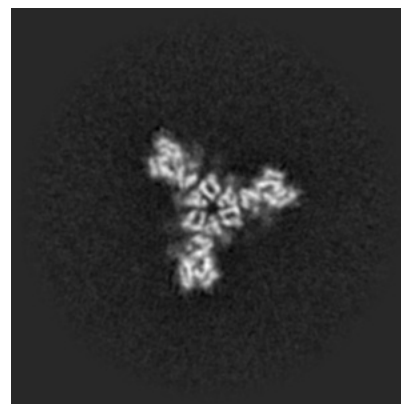
6.3.1 Primary map



X Index: 218



Y Index: 231



Z Index: 254

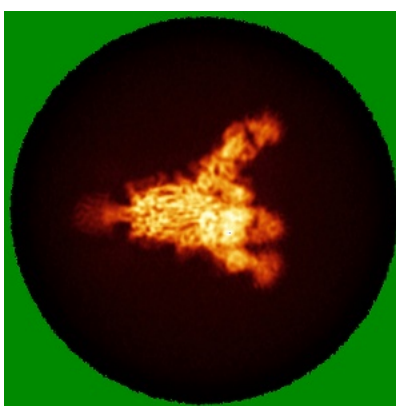
The images above show the largest variance slices of the map in three orthogonal directions.

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

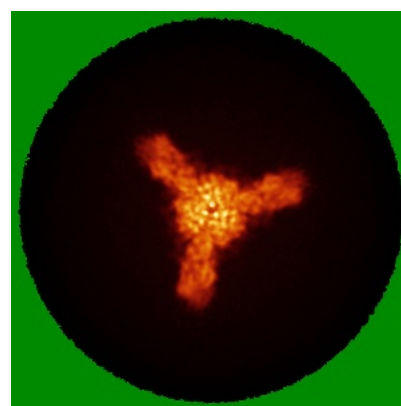
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.15. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

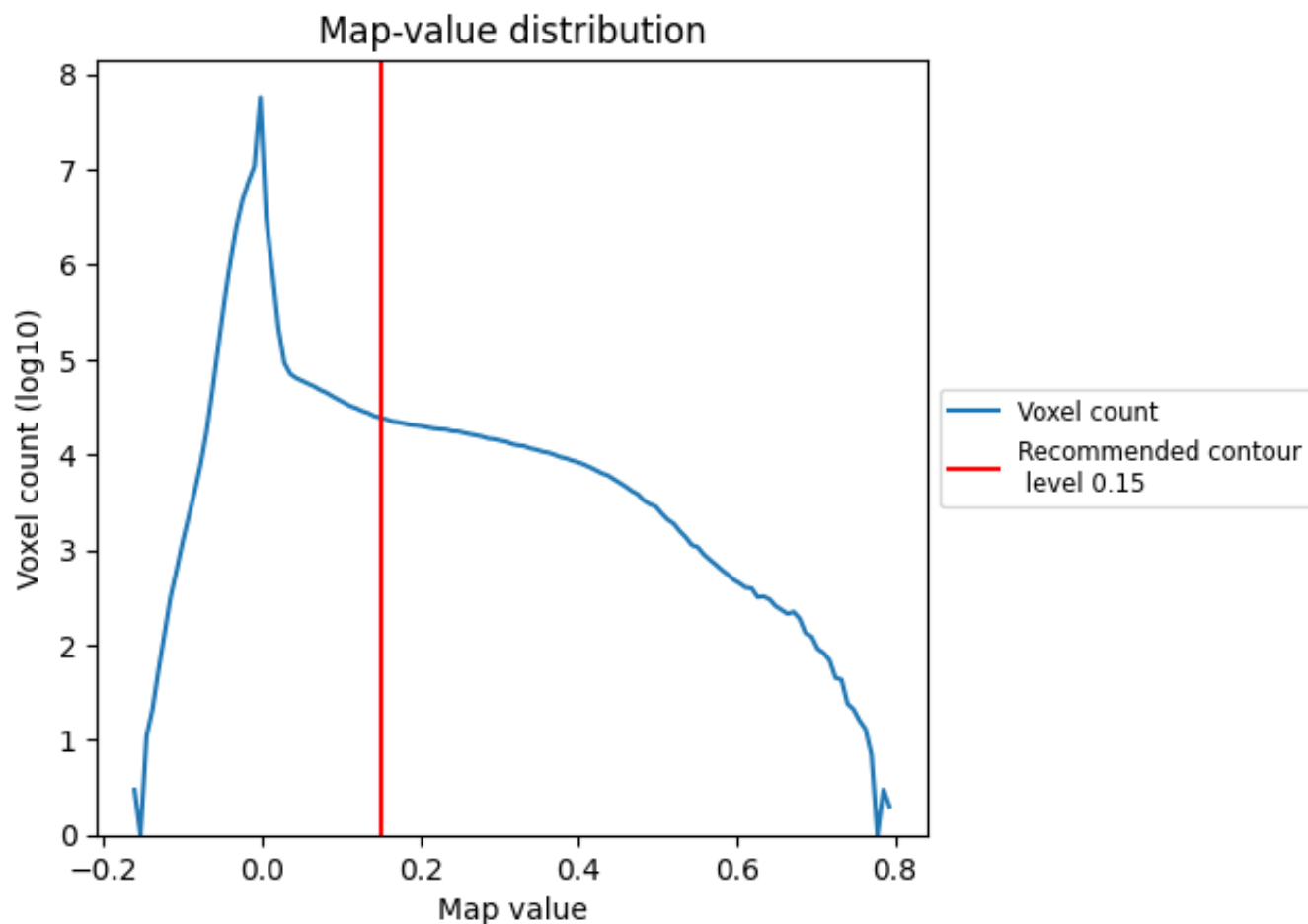
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

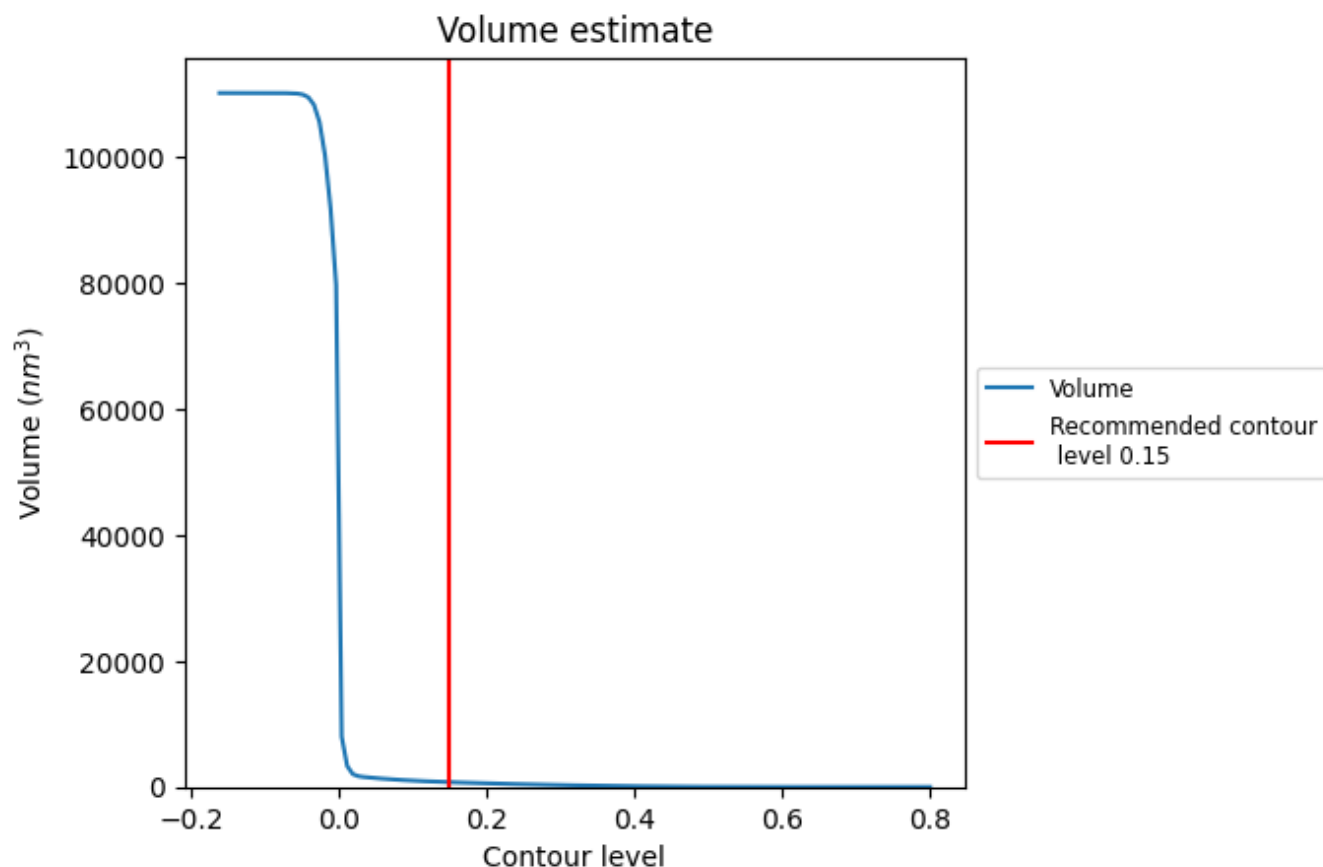
This section contains the results of statistical analysis of the map.

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.

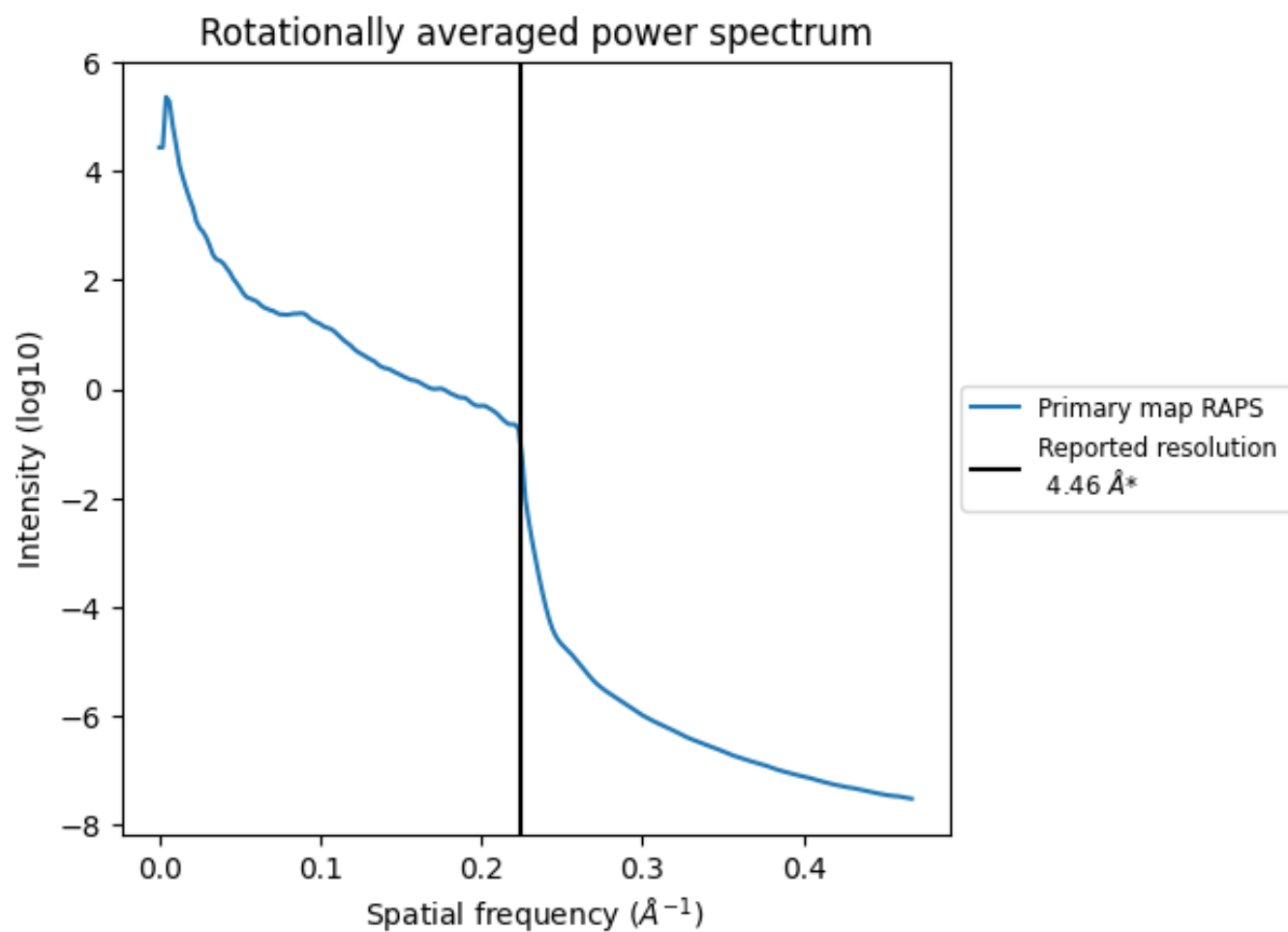
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 759 nm^3 ; this corresponds to an approximate mass of 686 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 0.224 Å⁻¹

8 Fourier-Shell correlation

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

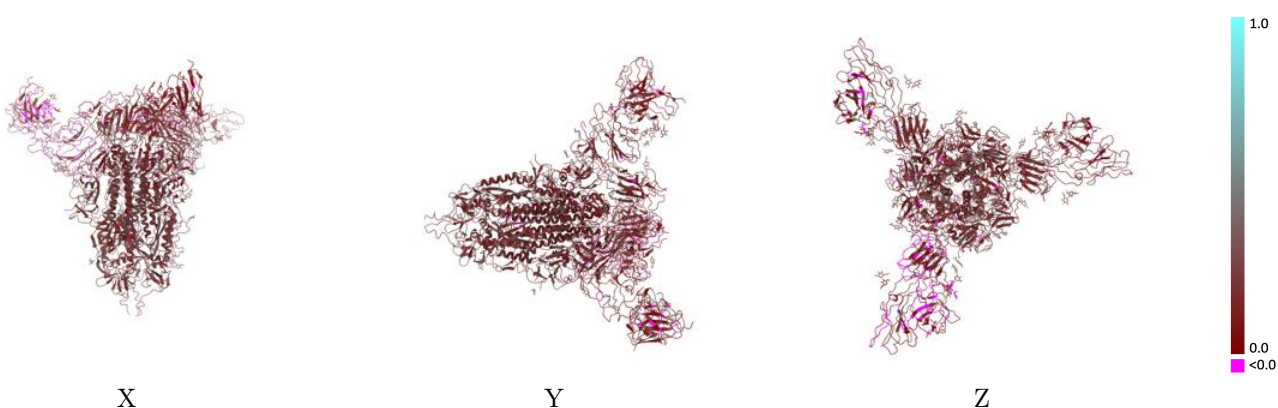
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-24075 and PDB model 7MXP. Per-residue inclusion information can be found in section 3 on page 14.

9.1 Map-model overlay [i](#)

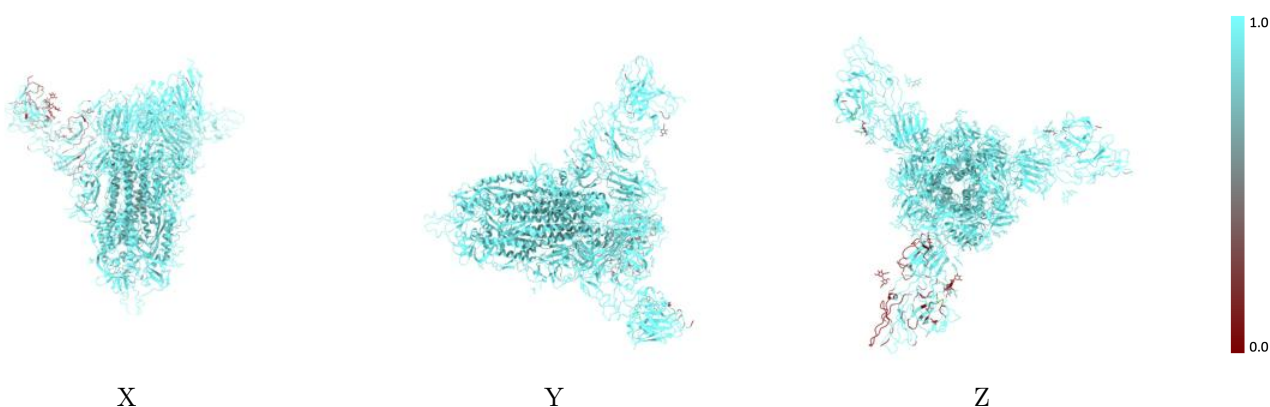
This section was not generated.

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



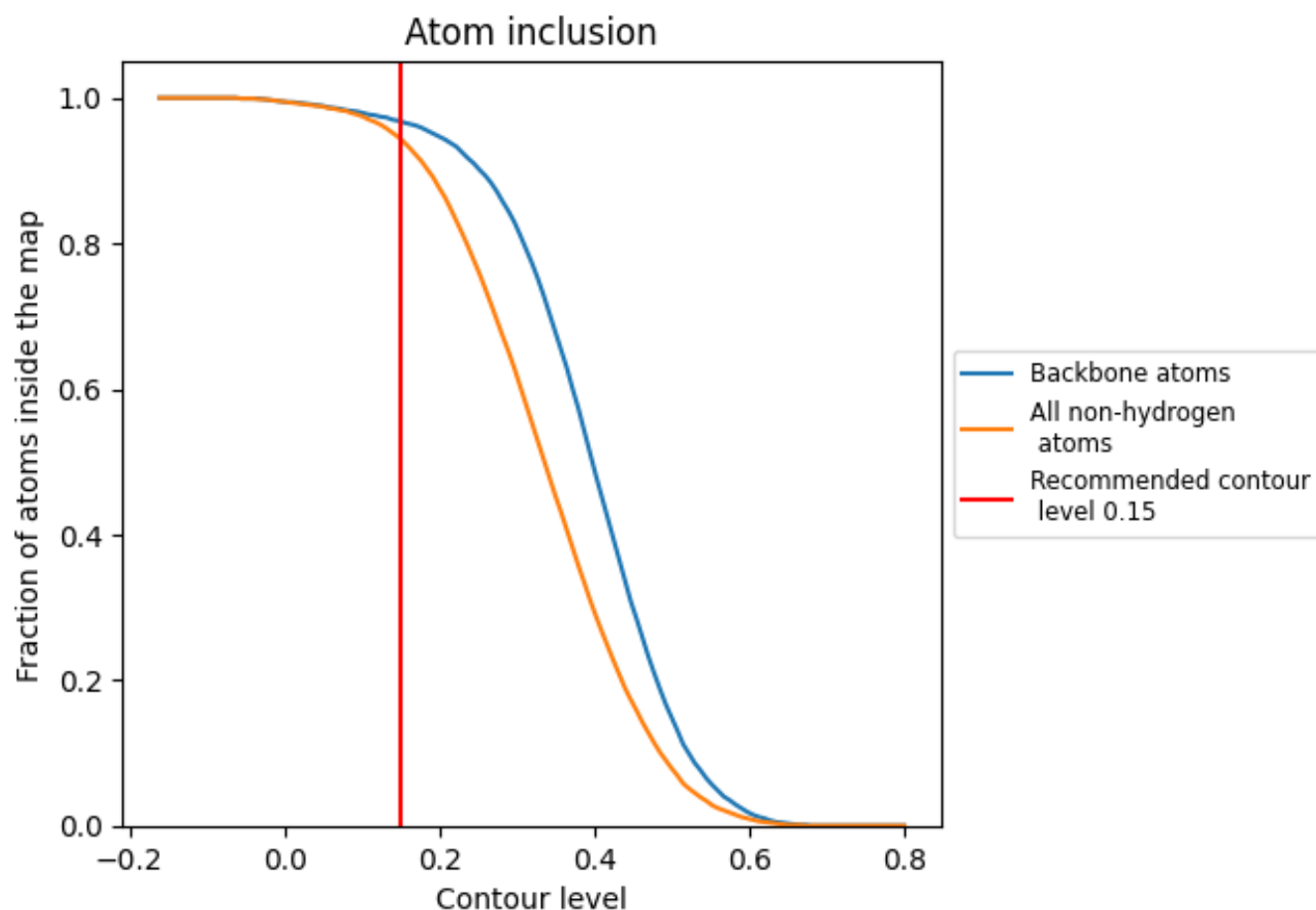
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.15).

9.4 Atom inclusion [i](#)



At the recommended contour level, 97% of all backbone atoms, 94% 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 (0.15) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9440	 0.1920
A	 0.9720	 0.2110
B	 0.9300	 0.1860
C	 0.9760	 0.2140
D	 0.0710	 0.1270
E	 0.3840	 0.0880
F	 0.7680	 0.2040
G	 0.7810	 0.0690
H	 0.9860	 0.1750
I	 0.9550	 0.1360
J	 0.6460	 0.0830
K	 0.9810	 0.1670
L	 0.9710	 0.1730
M	 0.8210	 0.2050
N	 0.8790	 0.2120
O	 0.9640	 0.1740
Y	 0.8970	 0.2380

