



Full wwPDB EM Validation Report ⓘ

Mar 20, 2026 – 02:30 AM UTC

PDB ID : 8Y5J / pdb_00008y5j
EMDB ID : EMD-38937
Title : Cryo-EM structure of SARS-CoV-2 Omicron JN.1 spike protein
Authors : Li, L.J.; Gu, Y.H.; Qi, J.X.; Gao, G.F.
Deposited on : 2024-01-31
Resolution : 2.94 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:

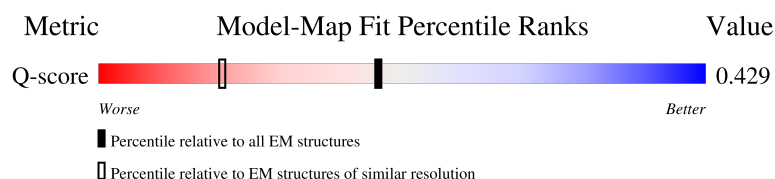
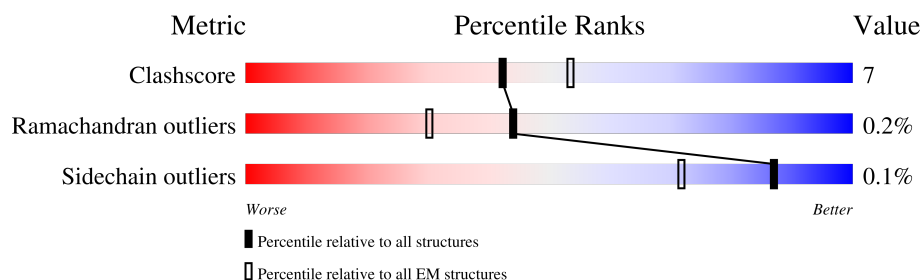
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 2.94 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	13068 (2.44 - 3.44)

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	1248	 10% 69% 12% 19%
1	B	1248	 10% 65% 14% 20%
1	C	1248	 10% 68% 13% 19%
2	D	2	 100%

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Mol	Chain	Length	Quality of chain
2	E	2	100%
2	F	2	50% 50%
3	G	3	100% 67% 33%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 24146 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	1011	Total	C	N	O	S	0	0
			7922	5073	1313	1500	36		
1	B	996	Total	C	N	O	S	0	0
			7815	5003	1294	1484	34		
1	C	1011	Total	C	N	O	S	0	0
			7922	5073	1313	1500	36		

There are 348 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	insertion	UNP P0DTC2
A	17	PRO	ASN	variant	UNP P0DTC2
A	19	PHE	THR	variant	UNP P0DTC2
A	20	ASN	THR	variant	UNP P0DTC2
A	21	LEU	ARG	variant	UNP P0DTC2
A	22	ILE	THR	variant	UNP P0DTC2
A	23	THR	GLN	variant	UNP P0DTC2
A	24	THR	LEU	variant	UNP P0DTC2
A	25	THR	PRO	variant	UNP P0DTC2
A	26	GLN	PRO	variant	UNP P0DTC2
A	27	SER	ALA	variant	UNP P0DTC2
A	50	LEU	SER	variant	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	127	PHE	VAL	variant	UNP P0DTC2
A	143	ASP	GLY	variant	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	157	SER	PHE	variant	UNP P0DTC2
A	158	GLY	ARG	variant	UNP P0DTC2
A	?	-	ASN	deletion	UNP P0DTC2
A	212	ILE	LEU	variant	UNP P0DTC2
A	213	GLY	VAL	variant	UNP P0DTC2
A	216	PHE	LEU	variant	UNP P0DTC2
A	245	ASN	HIS	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	264	ASP	ALA	variant	UNP P0DTC2
A	332	VAL	ILE	variant	UNP P0DTC2
A	339	HIS	GLY	variant	UNP P0DTC2
A	356	THR	LYS	variant	UNP P0DTC2
A	371	PHE	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	376	ALA	THR	variant	UNP P0DTC2
A	403	LYS	ARG	variant	UNP P0DTC2
A	405	ASN	ASP	variant	UNP P0DTC2
A	408	SER	ARG	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	445	HIS	VAL	variant	UNP P0DTC2
A	446	SER	GLY	variant	UNP P0DTC2
A	450	ASP	ASN	variant	UNP P0DTC2
A	452	TRP	LEU	variant	UNP P0DTC2
A	455	SER	LEU	variant	UNP P0DTC2
A	460	LYS	ASN	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	481	LYS	ASN	variant	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	483	LYS	GLU	variant	UNP P0DTC2
A	485	PRO	PHE	variant	UNP P0DTC2
A	497	ARG	GLN	variant	UNP P0DTC2
A	500	TYR	ASN	variant	UNP P0DTC2
A	504	HIS	TYR	variant	UNP P0DTC2
A	553	LYS	GLU	variant	UNP P0DTC2
A	569	VAL	ALA	variant	UNP P0DTC2
A	613	GLY	ASP	variant	UNP P0DTC2
A	620	SER	PRO	variant	UNP P0DTC2
A	654	TYR	HIS	variant	UNP P0DTC2
A	669	VAL	ILE	variant	UNP P0DTC2
A	678	LYS	ASN	variant	UNP P0DTC2
A	680	ARG	PRO	variant	UNP P0DTC2
A	763	LYS	ASN	variant	UNP P0DTC2
A	795	TYR	ASP	variant	UNP P0DTC2
A	816	PRO	PHE	engineered mutation	UNP P0DTC2
A	891	PRO	ALA	engineered mutation	UNP P0DTC2
A	898	PRO	ALA	engineered mutation	UNP P0DTC2
A	938	PHE	SER	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	941	PRO	ALA	engineered mutation	UNP P0DTC2
A	953	HIS	GLN	variant	UNP P0DTC2
A	968	LYS	ASN	variant	UNP P0DTC2
A	985	PRO	LYS	engineered mutation	UNP P0DTC2
A	986	PRO	VAL	engineered mutation	UNP P0DTC2
A	1142	LEU	PRO	variant	UNP P0DTC2
A	1208	GLY	-	expression tag	UNP P0DTC2
A	1209	GLY	-	expression tag	UNP P0DTC2
A	1210	GLY	-	expression tag	UNP P0DTC2
A	1211	SER	-	expression tag	UNP P0DTC2
A	1212	GLY	-	expression tag	UNP P0DTC2
A	1213	TYR	-	expression tag	UNP P0DTC2
A	1214	ILE	-	expression tag	UNP P0DTC2
A	1215	PRO	-	expression tag	UNP P0DTC2
A	1216	GLU	-	expression tag	UNP P0DTC2
A	1217	ALA	-	expression tag	UNP P0DTC2
A	1218	PRO	-	expression tag	UNP P0DTC2
A	1219	ARG	-	expression tag	UNP P0DTC2
A	1220	ASP	-	expression tag	UNP P0DTC2
A	1221	GLY	-	expression tag	UNP P0DTC2
A	1222	GLN	-	expression tag	UNP P0DTC2
A	1223	ALA	-	expression tag	UNP P0DTC2
A	1224	TYR	-	expression tag	UNP P0DTC2
A	1225	VAL	-	expression tag	UNP P0DTC2
A	1226	ARG	-	expression tag	UNP P0DTC2
A	1227	LYS	-	expression tag	UNP P0DTC2
A	1228	ASP	-	expression tag	UNP P0DTC2
A	1229	GLY	-	expression tag	UNP P0DTC2
A	1230	GLU	-	expression tag	UNP P0DTC2
A	1231	TRP	-	expression tag	UNP P0DTC2
A	1232	VAL	-	expression tag	UNP P0DTC2
A	1233	LEU	-	expression tag	UNP P0DTC2
A	1234	LEU	-	expression tag	UNP P0DTC2
A	1235	SER	-	expression tag	UNP P0DTC2
A	1236	THR	-	expression tag	UNP P0DTC2
A	1237	PHE	-	expression tag	UNP P0DTC2
A	1238	LEU	-	expression tag	UNP P0DTC2
A	1239	GLY	-	expression tag	UNP P0DTC2
A	1240	GLY	-	expression tag	UNP P0DTC2
A	1241	GLY	-	expression tag	UNP P0DTC2
A	1242	SER	-	expression tag	UNP P0DTC2
A	1243	ALA	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1244	TRP	-	expression tag	UNP P0DTC2
A	1245	SER	-	expression tag	UNP P0DTC2
A	1246	HIS	-	expression tag	UNP P0DTC2
A	1247	PRO	-	expression tag	UNP P0DTC2
A	1248	GLN	-	expression tag	UNP P0DTC2
A	1249	PHE	-	expression tag	UNP P0DTC2
A	1250	GLU	-	expression tag	UNP P0DTC2
A	1251	LYS	-	expression tag	UNP P0DTC2
B	16	MET	-	insertion	UNP P0DTC2
B	17	PRO	ASN	variant	UNP P0DTC2
B	19	PHE	THR	variant	UNP P0DTC2
B	20	ASN	THR	variant	UNP P0DTC2
B	21	LEU	ARG	variant	UNP P0DTC2
B	22	ILE	THR	variant	UNP P0DTC2
B	23	THR	GLN	variant	UNP P0DTC2
B	24	THR	LEU	variant	UNP P0DTC2
B	25	THR	PRO	variant	UNP P0DTC2
B	26	GLN	PRO	variant	UNP P0DTC2
B	27	SER	ALA	variant	UNP P0DTC2
B	50	LEU	SER	variant	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	127	PHE	VAL	variant	UNP P0DTC2
B	143	ASP	GLY	variant	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	157	SER	PHE	variant	UNP P0DTC2
B	158	GLY	ARG	variant	UNP P0DTC2
B	?	-	ASN	deletion	UNP P0DTC2
B	212	ILE	LEU	variant	UNP P0DTC2
B	213	GLY	VAL	variant	UNP P0DTC2
B	216	PHE	LEU	variant	UNP P0DTC2
B	245	ASN	HIS	variant	UNP P0DTC2
B	264	ASP	ALA	variant	UNP P0DTC2
B	332	VAL	ILE	variant	UNP P0DTC2
B	339	HIS	GLY	variant	UNP P0DTC2
B	356	THR	LYS	variant	UNP P0DTC2
B	371	PHE	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	376	ALA	THR	variant	UNP P0DTC2
B	403	LYS	ARG	variant	UNP P0DTC2
B	405	ASN	ASP	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	408	SER	ARG	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	445	HIS	VAL	variant	UNP P0DTC2
B	446	SER	GLY	variant	UNP P0DTC2
B	450	ASP	ASN	variant	UNP P0DTC2
B	452	TRP	LEU	variant	UNP P0DTC2
B	455	SER	LEU	variant	UNP P0DTC2
B	460	LYS	ASN	variant	UNP P0DTC2
B	477	ASN	SER	variant	UNP P0DTC2
B	478	LYS	THR	variant	UNP P0DTC2
B	481	LYS	ASN	variant	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	483	LYS	GLU	variant	UNP P0DTC2
B	485	PRO	PHE	variant	UNP P0DTC2
B	497	ARG	GLN	variant	UNP P0DTC2
B	500	TYR	ASN	variant	UNP P0DTC2
B	504	HIS	TYR	variant	UNP P0DTC2
B	553	LYS	GLU	variant	UNP P0DTC2
B	569	VAL	ALA	variant	UNP P0DTC2
B	613	GLY	ASP	variant	UNP P0DTC2
B	620	SER	PRO	variant	UNP P0DTC2
B	654	TYR	HIS	variant	UNP P0DTC2
B	669	VAL	ILE	variant	UNP P0DTC2
B	678	LYS	ASN	variant	UNP P0DTC2
B	680	ARG	PRO	variant	UNP P0DTC2
B	763	LYS	ASN	variant	UNP P0DTC2
B	795	TYR	ASP	variant	UNP P0DTC2
B	816	PRO	PHE	engineered mutation	UNP P0DTC2
B	891	PRO	ALA	engineered mutation	UNP P0DTC2
B	898	PRO	ALA	engineered mutation	UNP P0DTC2
B	938	PHE	SER	variant	UNP P0DTC2
B	941	PRO	ALA	engineered mutation	UNP P0DTC2
B	953	HIS	GLN	variant	UNP P0DTC2
B	968	LYS	ASN	variant	UNP P0DTC2
B	985	PRO	LYS	engineered mutation	UNP P0DTC2
B	986	PRO	VAL	engineered mutation	UNP P0DTC2
B	1142	LEU	PRO	variant	UNP P0DTC2
B	1208	GLY	-	expression tag	UNP P0DTC2
B	1209	GLY	-	expression tag	UNP P0DTC2
B	1210	GLY	-	expression tag	UNP P0DTC2
B	1211	SER	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1212	GLY	-	expression tag	UNP P0DTC2
B	1213	TYR	-	expression tag	UNP P0DTC2
B	1214	ILE	-	expression tag	UNP P0DTC2
B	1215	PRO	-	expression tag	UNP P0DTC2
B	1216	GLU	-	expression tag	UNP P0DTC2
B	1217	ALA	-	expression tag	UNP P0DTC2
B	1218	PRO	-	expression tag	UNP P0DTC2
B	1219	ARG	-	expression tag	UNP P0DTC2
B	1220	ASP	-	expression tag	UNP P0DTC2
B	1221	GLY	-	expression tag	UNP P0DTC2
B	1222	GLN	-	expression tag	UNP P0DTC2
B	1223	ALA	-	expression tag	UNP P0DTC2
B	1224	TYR	-	expression tag	UNP P0DTC2
B	1225	VAL	-	expression tag	UNP P0DTC2
B	1226	ARG	-	expression tag	UNP P0DTC2
B	1227	LYS	-	expression tag	UNP P0DTC2
B	1228	ASP	-	expression tag	UNP P0DTC2
B	1229	GLY	-	expression tag	UNP P0DTC2
B	1230	GLU	-	expression tag	UNP P0DTC2
B	1231	TRP	-	expression tag	UNP P0DTC2
B	1232	VAL	-	expression tag	UNP P0DTC2
B	1233	LEU	-	expression tag	UNP P0DTC2
B	1234	LEU	-	expression tag	UNP P0DTC2
B	1235	SER	-	expression tag	UNP P0DTC2
B	1236	THR	-	expression tag	UNP P0DTC2
B	1237	PHE	-	expression tag	UNP P0DTC2
B	1238	LEU	-	expression tag	UNP P0DTC2
B	1239	GLY	-	expression tag	UNP P0DTC2
B	1240	GLY	-	expression tag	UNP P0DTC2
B	1241	GLY	-	expression tag	UNP P0DTC2
B	1242	SER	-	expression tag	UNP P0DTC2
B	1243	ALA	-	expression tag	UNP P0DTC2
B	1244	TRP	-	expression tag	UNP P0DTC2
B	1245	SER	-	expression tag	UNP P0DTC2
B	1246	HIS	-	expression tag	UNP P0DTC2
B	1247	PRO	-	expression tag	UNP P0DTC2
B	1248	GLN	-	expression tag	UNP P0DTC2
B	1249	PHE	-	expression tag	UNP P0DTC2
B	1250	GLU	-	expression tag	UNP P0DTC2
B	1251	LYS	-	expression tag	UNP P0DTC2
C	16	MET	-	insertion	UNP P0DTC2
C	17	PRO	ASN	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	19	PHE	THR	variant	UNP P0DTC2
C	20	ASN	THR	variant	UNP P0DTC2
C	21	LEU	ARG	variant	UNP P0DTC2
C	22	ILE	THR	variant	UNP P0DTC2
C	23	THR	GLN	variant	UNP P0DTC2
C	24	THR	LEU	variant	UNP P0DTC2
C	25	THR	PRO	variant	UNP P0DTC2
C	26	GLN	PRO	variant	UNP P0DTC2
C	27	SER	ALA	variant	UNP P0DTC2
C	50	LEU	SER	variant	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	127	PHE	VAL	variant	UNP P0DTC2
C	143	ASP	GLY	variant	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	157	SER	PHE	variant	UNP P0DTC2
C	158	GLY	ARG	variant	UNP P0DTC2
C	?	-	ASN	deletion	UNP P0DTC2
C	212	ILE	LEU	variant	UNP P0DTC2
C	213	GLY	VAL	variant	UNP P0DTC2
C	216	PHE	LEU	variant	UNP P0DTC2
C	245	ASN	HIS	variant	UNP P0DTC2
C	264	ASP	ALA	variant	UNP P0DTC2
C	332	VAL	ILE	variant	UNP P0DTC2
C	339	HIS	GLY	variant	UNP P0DTC2
C	356	THR	LYS	variant	UNP P0DTC2
C	371	PHE	SER	variant	UNP P0DTC2
C	373	PRO	SER	variant	UNP P0DTC2
C	375	PHE	SER	variant	UNP P0DTC2
C	376	ALA	THR	variant	UNP P0DTC2
C	403	LYS	ARG	variant	UNP P0DTC2
C	405	ASN	ASP	variant	UNP P0DTC2
C	408	SER	ARG	variant	UNP P0DTC2
C	417	ASN	LYS	variant	UNP P0DTC2
C	440	LYS	ASN	variant	UNP P0DTC2
C	445	HIS	VAL	variant	UNP P0DTC2
C	446	SER	GLY	variant	UNP P0DTC2
C	450	ASP	ASN	variant	UNP P0DTC2
C	452	TRP	LEU	variant	UNP P0DTC2
C	455	SER	LEU	variant	UNP P0DTC2
C	460	LYS	ASN	variant	UNP P0DTC2
C	477	ASN	SER	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	478	LYS	THR	variant	UNP P0DTC2
C	481	LYS	ASN	variant	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	483	LYS	GLU	variant	UNP P0DTC2
C	485	PRO	PHE	variant	UNP P0DTC2
C	497	ARG	GLN	variant	UNP P0DTC2
C	500	TYR	ASN	variant	UNP P0DTC2
C	504	HIS	TYR	variant	UNP P0DTC2
C	553	LYS	GLU	variant	UNP P0DTC2
C	569	VAL	ALA	variant	UNP P0DTC2
C	613	GLY	ASP	variant	UNP P0DTC2
C	620	SER	PRO	variant	UNP P0DTC2
C	654	TYR	HIS	variant	UNP P0DTC2
C	669	VAL	ILE	variant	UNP P0DTC2
C	678	LYS	ASN	variant	UNP P0DTC2
C	680	ARG	PRO	variant	UNP P0DTC2
C	763	LYS	ASN	variant	UNP P0DTC2
C	795	TYR	ASP	variant	UNP P0DTC2
C	816	PRO	PHE	engineered mutation	UNP P0DTC2
C	891	PRO	ALA	engineered mutation	UNP P0DTC2
C	898	PRO	ALA	engineered mutation	UNP P0DTC2
C	938	PHE	SER	variant	UNP P0DTC2
C	941	PRO	ALA	engineered mutation	UNP P0DTC2
C	953	HIS	GLN	variant	UNP P0DTC2
C	968	LYS	ASN	variant	UNP P0DTC2
C	985	PRO	LYS	engineered mutation	UNP P0DTC2
C	986	PRO	VAL	engineered mutation	UNP P0DTC2
C	1142	LEU	PRO	variant	UNP P0DTC2
C	1208	GLY	-	expression tag	UNP P0DTC2
C	1209	GLY	-	expression tag	UNP P0DTC2
C	1210	GLY	-	expression tag	UNP P0DTC2
C	1211	SER	-	expression tag	UNP P0DTC2
C	1212	GLY	-	expression tag	UNP P0DTC2
C	1213	TYR	-	expression tag	UNP P0DTC2
C	1214	ILE	-	expression tag	UNP P0DTC2
C	1215	PRO	-	expression tag	UNP P0DTC2
C	1216	GLU	-	expression tag	UNP P0DTC2
C	1217	ALA	-	expression tag	UNP P0DTC2
C	1218	PRO	-	expression tag	UNP P0DTC2
C	1219	ARG	-	expression tag	UNP P0DTC2
C	1220	ASP	-	expression tag	UNP P0DTC2
C	1221	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1222	GLN	-	expression tag	UNP P0DTC2
C	1223	ALA	-	expression tag	UNP P0DTC2
C	1224	TYR	-	expression tag	UNP P0DTC2
C	1225	VAL	-	expression tag	UNP P0DTC2
C	1226	ARG	-	expression tag	UNP P0DTC2
C	1227	LYS	-	expression tag	UNP P0DTC2
C	1228	ASP	-	expression tag	UNP P0DTC2
C	1229	GLY	-	expression tag	UNP P0DTC2
C	1230	GLU	-	expression tag	UNP P0DTC2
C	1231	TRP	-	expression tag	UNP P0DTC2
C	1232	VAL	-	expression tag	UNP P0DTC2
C	1233	LEU	-	expression tag	UNP P0DTC2
C	1234	LEU	-	expression tag	UNP P0DTC2
C	1235	SER	-	expression tag	UNP P0DTC2
C	1236	THR	-	expression tag	UNP P0DTC2
C	1237	PHE	-	expression tag	UNP P0DTC2
C	1238	LEU	-	expression tag	UNP P0DTC2
C	1239	GLY	-	expression tag	UNP P0DTC2
C	1240	GLY	-	expression tag	UNP P0DTC2
C	1241	GLY	-	expression tag	UNP P0DTC2
C	1242	SER	-	expression tag	UNP P0DTC2
C	1243	ALA	-	expression tag	UNP P0DTC2
C	1244	TRP	-	expression tag	UNP P0DTC2
C	1245	SER	-	expression tag	UNP P0DTC2
C	1246	HIS	-	expression tag	UNP P0DTC2
C	1247	PRO	-	expression tag	UNP P0DTC2
C	1248	GLN	-	expression tag	UNP P0DTC2
C	1249	PHE	-	expression tag	UNP P0DTC2
C	1250	GLU	-	expression tag	UNP P0DTC2
C	1251	LYS	-	expression tag	UNP P0DTC2

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



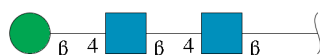
Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	2	Total	C	N	O	0	0
			28	16	2	10		

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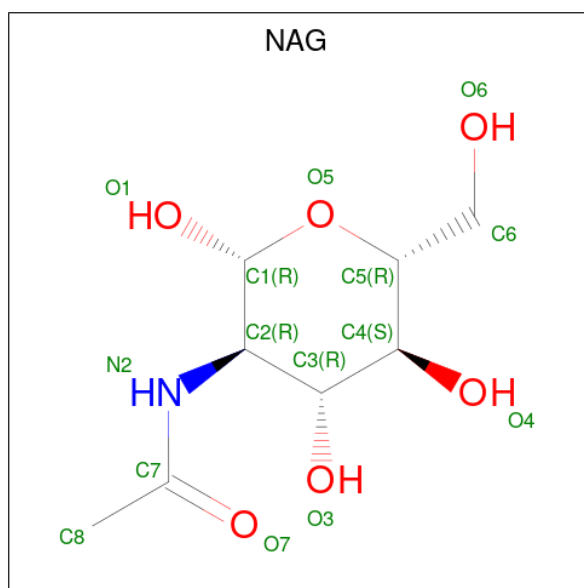
Mol	Chain	Residues	Atoms				AltConf	Trace
2	E	2	Total	C	N	O	0	0
			28	16	2	10		
2	F	2	Total	C	N	O	0	0
			28	16	2	10		

- 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	G	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	

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

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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 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	378152	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.974	Depositor
Minimum map value	-0.320	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.026	Depositor
Recommended contour level	0.2	Depositor
Map size ($\text{\AA}$)	391.0, 391.0, 391.0	wwPDB
Map dimensions	460, 460, 460	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.85, 0.85, 0.85	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.17	0/8115	0.34	0/11046
1	B	0.19	0/8003	0.39	0/10891
1	C	0.18	0/8115	0.35	0/11046
All	All	0.18	0/24233	0.36	0/32983

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	7922	0	7723	110	0
1	B	7815	0	7605	132	0
1	C	7922	0	7722	99	0
2	D	28	0	25	0	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
3	G	39	0	34	0	0
4	A	140	0	130	3	0
4	B	112	0	104	2	0
4	C	112	0	104	1	0
All	All	24146	0	23497	320	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 7.

All (320) 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:329:PHE:CE2	1:A:527:LYS:HB2	1.77	1.18
1:A:329:PHE:CE2	1:A:527:LYS:HD2	1.83	1.14
1:B:431:GLY:HA3	1:B:512:LEU:O	1.61	1.00
1:A:329:PHE:HE2	1:A:527:LYS:HD2	1.30	0.96
1:A:329:PHE:CD2	1:A:527:LYS:HB2	2.03	0.93
1:A:329:PHE:CZ	1:A:527:LYS:HD2	2.04	0.91
1:B:334:ASN:HB2	1:B:361:CYS:HA	1.52	0.90
1:A:329:PHE:HE2	1:A:527:LYS:CD	1.86	0.89
1:A:329:PHE:CE2	1:A:527:LYS:CB	2.56	0.88
1:A:329:PHE:HE2	1:A:527:LYS:CB	1.89	0.83
1:A:329:PHE:HE2	1:A:527:LYS:HB2	1.37	0.80
1:A:329:PHE:CE2	1:A:527:LYS:CD	2.61	0.80
1:B:131:CYS:HA	1:B:166:CYS:HB3	1.64	0.80
1:A:329:PHE:HE2	1:A:527:LYS:CG	1.94	0.79
1:B:334:ASN:CB	1:B:361:CYS:HA	2.13	0.79
1:B:332:VAL:HG22	1:B:332:VAL:O	1.81	0.78
1:B:358:ILE:H	1:B:395:VAL:HG11	1.49	0.74
1:C:384:PRO:HD2	1:C:387:LEU:HD12	1.70	0.74
1:A:393:THR:HG21	1:A:518:HIS:HB2	1.71	0.71
1:B:400:PHE:HB2	1:B:509:VAL:HB	1.72	0.71
1:B:452:TRP:HB3	1:B:491:LEU:HB3	1.72	0.70
1:B:119:ILE:HG12	1:B:128:ILE:HG12	1.73	0.70
1:A:390:LEU:HB2	1:A:544:GLY:HA3	1.74	0.69
1:C:125:ASN:HB2	1:C:172:SER:HA	1.74	0.68
1:A:327:VAL:CG1	1:A:329:PHE:CZ	2.78	0.67
1:A:320:VAL:HG23	1:A:627:GLN:HB3	1.75	0.67
1:A:675:THR:HA	1:A:689:GLN:HG3	1.77	0.67
1:B:92:PHE:HE1	1:B:94:SER:HB3	1.60	0.66
1:A:391:CYS:N	1:A:524:CYS:SG	2.67	0.66
1:B:126:VAL:H	1:B:172:SER:HB3	1.61	0.65
1:B:824:LYS:NZ	1:B:937:LEU:O	2.25	0.65
1:B:904:ARG:NH1	1:B:1048:LEU:O	2.30	0.64
1:C:376:ALA:HB3	1:C:435:ALA:HB3	1.79	0.64
1:A:488:TYR:OH	1:B:375:PHE:HA	1.97	0.64
1:C:301:CYS:O	1:C:304:LYS:NZ	2.31	0.63
1:A:758:PHE:HD2	1:A:1000:LEU:HD11	1.62	0.63
1:C:106:PHE:HB2	1:C:117:LEU:HB3	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:904:ARG:NH1	1:A:1048:LEU:O	2.31	0.63
1:C:444:LYS:HE3	1:C:448:ASN:HA	1.80	0.62
1:C:904:ARG:NH1	1:C:1048:LEU:O	2.31	0.62
1:B:376:ALA:HB3	1:B:435:ALA:HB3	1.81	0.62
1:A:698:LEU:HB3	1:B:872:TYR:HE1	1.66	0.60
1:B:231:ILE:HD12	1:B:233:ILE:HG12	1.83	0.60
1:B:395:VAL:HG23	1:B:514:PHE:HE1	1.66	0.60
1:B:320:VAL:HG23	1:B:627:GLN:HB3	1.83	0.60
1:C:382:VAL:HG22	1:C:384:PRO:HD3	1.82	0.60
1:A:872:TYR:HE1	1:C:698:LEU:HB3	1.66	0.60
1:C:357:ARG:NH1	1:C:394:ASN:OD1	2.35	0.60
1:B:356:THR:HB	1:B:397:ALA:HB3	1.85	0.59
1:B:417:ASN:O	1:B:421:TYR:HB2	2.03	0.59
1:A:403:LYS:HG2	1:A:496:PHE:HE1	1.68	0.59
1:A:374:PHE:HB3	1:A:436:TRP:HB3	1.84	0.59
1:B:616:CYS:SG	1:B:643:GLN:NE2	2.76	0.58
1:C:641:VAL:HG13	1:C:650:ILE:HG12	1.85	0.58
1:A:133:PHE:HA	1:A:162:SER:HB3	1.85	0.58
1:A:643:GLN:NE2	1:A:644:THR:O	2.36	0.58
1:B:357:ARG:NH2	1:B:393:THR:O	2.36	0.58
1:B:431:GLY:CA	1:B:512:LEU:O	2.44	0.58
1:C:381:GLY:HA3	1:C:430:THR:HA	1.84	0.58
1:B:472:ILE:HG22	1:B:490:PRO:HD3	1.84	0.58
1:B:762:LEU:HD22	1:B:1007:VAL:HG21	1.86	0.58
1:C:522:THR:HG23	1:C:523:VAL:HG23	1.84	0.58
1:A:658:SER:HB3	1:A:697:SER:HB3	1.86	0.58
1:B:1073:ASN:OD1	1:C:894:GLN:NE2	2.36	0.57
1:B:337:PRO:HG2	1:B:358:ILE:HD13	1.86	0.57
1:B:193:VAL:HG23	1:B:223:LEU:HD22	1.86	0.57
1:B:327:VAL:HG13	1:B:541:ASN:HB3	1.85	0.57
1:B:392:PHE:HE2	1:B:395:VAL:HB	1.68	0.57
1:B:776:ASN:OD1	1:B:1018:ARG:NH1	2.34	0.57
1:A:457:ARG:HG2	1:A:459:SER:H	1.69	0.57
1:B:206:LYS:NZ	1:B:222:ALA:O	2.35	0.57
1:B:133:PHE:HB2	1:B:161:SER:HB3	1.86	0.57
1:B:137:ASN:ND2	1:B:157:SER:O	2.38	0.56
1:A:421:TYR:HA	1:A:461:LEU:HD23	1.87	0.56
1:B:334:ASN:HB2	1:B:361:CYS:CA	2.31	0.56
1:C:290:ASP:OD1	1:C:291:CYS:N	2.38	0.56
1:A:604:SER:OG	1:A:605:ASN:N	2.39	0.56
1:B:206:LYS:HB2	1:B:223:LEU:HG	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:132:GLU:HB3	1:C:165:ASN:HB2	1.87	0.55
1:A:824:LYS:NZ	1:A:937:LEU:O	2.28	0.55
1:A:92:PHE:HE1	1:A:94:SER:HB3	1.72	0.54
1:B:379:CYS:HB3	1:B:382:VAL:O	2.08	0.54
1:B:758:PHE:HD2	1:B:1000:LEU:HD11	1.73	0.54
1:C:323:THR:OG1	1:C:324:GLU:OE1	2.22	0.54
1:A:454:ARG:HD3	1:A:490:PRO:HB2	1.89	0.54
1:B:911:THR:OG1	1:B:1105:GLN:NE2	2.39	0.54
1:B:362:VAL:HG13	1:B:388:ASN:HA	1.90	0.54
1:A:356:THR:HG23	4:A:1310:NAG:H62	1.89	0.53
1:C:658:SER:HB3	1:C:697:SER:HB3	1.89	0.53
1:B:128:ILE:HB	1:B:170:TYR:HB3	1.91	0.53
1:C:604:SER:OG	1:C:605:ASN:N	2.40	0.53
1:A:110:LEU:HD22	1:A:239:GLN:HE21	1.73	0.53
1:B:93:ALA:HA	1:B:191:GLU:HA	1.91	0.53
1:C:758:PHE:HD2	1:C:1000:LEU:HD11	1.74	0.53
1:B:732:LYS:HD2	1:B:770:ALA:HB1	1.90	0.52
1:A:190:ARG:HB3	1:A:192:PHE:HE1	1.74	0.52
1:B:641:VAL:HG22	1:B:650:ILE:HG23	1.92	0.52
1:C:368:LEU:HA	1:C:371:PHE:HD2	1.74	0.52
1:A:422:ASN:HD21	1:A:453:TYR:HD2	1.57	0.52
1:C:641:VAL:HG22	1:C:650:ILE:HG23	1.91	0.52
1:B:500:TYR:HB3	1:B:504:HIS:HB2	1.92	0.52
1:C:597:ILE:HB	1:C:608:ALA:HB3	1.90	0.52
1:A:421:TYR:CD1	1:A:457:ARG:HB3	2.45	0.52
1:C:46:SER:OG	1:C:281:GLU:HG2	2.10	0.52
1:B:442:ASP:CG	1:B:508:ARG:HH21	2.19	0.51
1:C:330:PRO:HB3	1:C:578:PRO:HB2	1.91	0.51
1:B:857:LEU:HD13	1:B:958:LEU:HD22	1.92	0.51
1:B:485:PRO:O	1:B:487:CYS:N	2.41	0.51
1:B:44:ARG:HB2	1:B:279:TYR:CD2	2.46	0.51
1:A:497:ARG:NH1	1:A:500:TYR:OH	2.43	0.51
1:B:332:VAL:HA	4:B:1308:NAG:H82	1.93	0.51
1:C:988:ALA:O	1:C:992:ILE:HG12	2.11	0.51
1:A:383:SER:H	1:B:983:LEU:HD23	1.76	0.51
1:B:333:THR:O	1:B:334:ASN:ND2	2.44	0.51
1:C:472:ILE:HG22	1:C:489:PHE:HD1	1.76	0.50
1:B:701:GLU:HB3	1:C:789:LYS:NZ	2.27	0.50
1:B:336:CYS:O	1:B:338:PHE:N	2.42	0.50
1:B:419:ALA:HB1	1:B:424:LYS:HD2	1.93	0.50
1:C:762:LEU:HD22	1:C:1007:VAL:HG21	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:282:ASN:O	1:C:284:THR:N	2.44	0.50
1:B:375:PHE:O	1:B:376:ALA:C	2.55	0.50
1:B:454:ARG:HD3	1:B:457:ARG:HB2	1.94	0.50
1:B:604:SER:OG	1:B:605:ASN:N	2.44	0.50
1:B:376:ALA:HA	1:B:378:LYS:HZ3	1.76	0.49
1:A:126:VAL:HG12	1:A:172:SER:HB3	1.93	0.49
1:B:820:LEU:O	1:B:824:LYS:HB2	2.11	0.49
1:A:698:LEU:HD22	1:B:872:TYR:CZ	2.47	0.49
1:B:106:PHE:CD1	1:B:238:PHE:HB2	2.47	0.49
1:C:422:ASN:HD21	1:C:453:TYR:HD2	1.60	0.49
1:A:290:ASP:OD1	1:A:291:CYS:N	2.45	0.49
1:C:755:TYR:OH	1:C:993:ASP:OD1	2.26	0.49
1:B:374:PHE:HD1	1:B:436:TRP:HB3	1.77	0.49
1:A:138:ASP:N	1:A:138:ASP:OD1	2.46	0.49
1:A:1027:LYS:NZ	1:A:1041:PHE:O	2.45	0.49
1:B:410:ILE:HG21	1:B:433:VAL:HG11	1.95	0.49
1:A:762:LEU:HD22	1:A:1007:VAL:HG21	1.95	0.49
1:C:562:GLN:O	1:C:576:ARG:NH2	2.46	0.49
1:C:675:THR:HG22	1:C:689:GLN:HG3	1.95	0.49
1:A:379:CYS:HA	1:A:432:CYS:HA	1.95	0.48
1:A:988:ALA:O	1:A:992:ILE:HG12	2.13	0.48
1:B:804:ILE:HD12	1:B:877:LEU:HD11	1.95	0.48
1:B:810:LYS:NZ	1:B:819:ASP:OD2	2.31	0.48
1:C:353:TRP:CD1	1:C:353:TRP:H	2.30	0.48
1:C:911:THR:OG1	1:C:1105:GLN:NE2	2.44	0.48
1:A:312:ILE:HD13	1:A:597:ILE:HG13	1.95	0.48
1:A:776:ASN:OD1	1:A:1018:ARG:NH1	2.47	0.48
1:A:791:PRO:HG2	1:C:706:TYR:HB3	1.95	0.48
1:C:878:ALA:O	1:C:882:THR:OG1	2.27	0.48
1:B:332:VAL:O	1:B:333:THR:C	2.56	0.48
1:C:905:PHE:CD2	1:C:915:LEU:HB2	2.49	0.48
1:B:356:THR:HG22	1:B:358:ILE:HD11	1.96	0.48
1:B:404:GLY:HA2	1:B:507:TYR:CE2	2.49	0.48
1:A:953:HIS:HB3	1:A:1013:ARG:CZ	2.43	0.47
1:B:421:TYR:HD1	1:B:457:ARG:HB3	1.80	0.47
1:C:130:VAL:HB	1:C:168:PHE:HB3	1.96	0.47
1:B:988:ALA:O	1:B:992:ILE:HG12	2.15	0.47
1:C:290:ASP:O	1:C:297:SER:OG	2.32	0.47
1:B:44:ARG:HB2	1:B:279:TYR:CE2	2.50	0.47
1:B:106:PHE:HD2	1:B:117:LEU:HD23	1.80	0.47
1:C:409:GLN:HA	1:C:414:GLN:HG2	1.94	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:LYS:HG3	1:A:208:THR:HG23	1.96	0.47
1:A:392:PHE:HB2	1:A:523:VAL:HG13	1.95	0.47
1:B:341:VAL:HG21	1:B:356:THR:HG23	1.95	0.47
1:B:403:LYS:HG3	1:B:494:TYR:CE1	2.50	0.47
1:B:434:ILE:O	1:B:509:VAL:HA	2.14	0.47
1:C:620:SER:O	1:C:624:HIS:ND1	2.36	0.47
1:B:193:VAL:HB	1:B:204:TYR:HB2	1.97	0.47
1:C:50:LEU:HD22	1:C:304:LYS:NZ	2.30	0.47
1:C:295:PRO:HG3	1:C:632:TRP:CE3	2.50	0.47
1:B:448:ASN:HB3	1:B:496:PHE:HB2	1.96	0.47
1:C:358:ILE:HB	1:C:395:VAL:HG11	1.97	0.47
1:C:1096:SER:HB2	1:C:1101:TRP:CD2	2.50	0.47
1:A:472:ILE:HG21	1:A:481:LYS:HG2	1.97	0.46
1:C:711:ILE:HD13	1:C:1093:VAL:HG11	1.96	0.46
1:C:732:LYS:HD2	1:C:770:ALA:HB1	1.97	0.46
1:C:725:ILE:HG13	1:C:1060:VAL:HG22	1.98	0.46
1:C:976:LEU:HD22	1:C:992:ILE:HD12	1.97	0.46
1:C:428:ASP:OD1	1:C:428:ASP:N	2.45	0.46
1:B:375:PHE:O	1:B:435:ALA:O	2.34	0.46
1:C:574:ALA:HB1	1:C:583:ILE:HD11	1.98	0.46
1:C:1081:CYS:HB2	1:C:1131:ILE:HD11	1.98	0.46
1:A:92:PHE:CE1	1:A:94:SER:HB3	2.51	0.45
1:C:206:LYS:HB2	1:C:223:LEU:HG	1.98	0.45
1:C:316:SER:O	1:C:594:VAL:HG12	2.16	0.45
1:A:920:LYS:HD2	1:A:920:LYS:HA	1.73	0.45
1:B:344:ALA:HB3	1:B:347:PHE:HE1	1.81	0.45
1:A:706:TYR:HB3	1:B:791:PRO:HG2	1.98	0.45
1:C:310:LYS:HG3	1:C:599:PRO:HA	1.99	0.45
1:C:553:LYS:NZ	1:C:554:SER:O	2.48	0.45
1:C:637:THR:H	1:C:640:ASN:HD21	1.64	0.45
1:B:985:PRO:N	1:B:986:PRO:HD2	2.32	0.45
1:C:128:ILE:HD13	1:C:229:LEU:HD11	1.98	0.45
1:B:641:VAL:HG13	1:B:650:ILE:HG12	1.97	0.45
1:B:290:ASP:OD1	1:B:291:CYS:N	2.46	0.45
1:A:327:VAL:HG12	1:A:329:PHE:CZ	2.51	0.45
1:B:331:ASN:HB2	1:B:579:GLN:HA	1.97	0.45
1:C:320:VAL:HG23	1:C:627:GLN:HB3	1.98	0.45
1:A:102:ARG:HA	1:A:102:ARG:HD3	1.74	0.45
1:A:364:ASP:OD1	1:A:366:SER:OG	2.32	0.45
1:A:698:LEU:HB3	1:B:872:TYR:CE1	2.50	0.45
1:A:136:CYS:HA	1:A:160:TYR:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:984:ASP:OD1	1:B:984:ASP:N	2.49	0.44
1:A:824:LYS:NZ	1:A:940:THR:O	2.33	0.44
1:A:872:TYR:CZ	1:C:698:LEU:HD22	2.52	0.44
1:B:421:TYR:O	1:B:454:ARG:HD2	2.17	0.44
1:A:885:TRP:H	1:A:885:TRP:HE3	1.65	0.44
1:B:755:TYR:OH	1:B:993:ASP:OD1	2.27	0.44
1:C:379:CYS:HB2	1:C:383:SER:HB3	1.99	0.44
1:C:452:TRP:CZ2	1:C:493:SER:HB3	2.52	0.44
1:A:452:TRP:CD1	1:A:493:SER:HA	2.53	0.44
1:A:483:LYS:HE3	1:B:375:PHE:CE1	2.52	0.44
1:B:130:VAL:O	1:B:130:VAL:HG13	2.17	0.44
1:C:987:GLU:O	1:C:991:GLN:HG2	2.17	0.44
1:A:872:TYR:OH	1:C:698:LEU:HD22	2.17	0.44
1:A:985:PRO:N	1:A:986:PRO:HD2	2.32	0.44
1:B:927:ASN:O	1:B:930:ILE:HG22	2.18	0.44
1:B:580:THR:O	1:B:581:LEU:C	2.61	0.44
1:A:199:GLY:HA2	1:A:232:GLY:HA2	1.99	0.44
1:A:295:PRO:HG3	1:A:632:TRP:CZ3	2.53	0.44
1:A:337:PRO:HG2	1:A:358:ILE:HG23	1.99	0.44
1:A:1100:HIS:ND1	4:A:1305:NAG:H5	2.33	0.44
1:B:398:ASP:HB2	1:B:511:VAL:HB	1.98	0.44
1:B:1027:LYS:NZ	1:B:1041:PHE:O	2.49	0.44
1:A:328:ARG:O	1:A:329:PHE:HD1	2.01	0.44
1:C:331:ASN:OD1	1:C:331:ASN:N	2.51	0.44
1:A:323:THR:OG1	1:A:324:GLU:OE1	2.34	0.44
1:A:328:ARG:NH1	1:A:530:THR:O	2.51	0.44
1:A:287:ASP:HB3	1:A:306:PHE:HE2	1.83	0.43
1:A:857:LEU:HD13	1:A:958:LEU:HD22	2.00	0.43
1:B:332:VAL:O	1:B:332:VAL:CG2	2.51	0.43
1:B:329:PHE:O	1:B:330:PRO:C	2.61	0.43
1:B:946:LYS:HB3	1:B:946:LYS:HE2	1.84	0.43
1:A:453:TYR:CE1	1:A:492:GLN:HB3	2.52	0.43
1:C:405:ASN:N	1:C:503:GLY:O	2.49	0.43
1:A:868:MET:HG2	1:C:698:LEU:HD21	2.00	0.43
1:B:104:TRP:N	1:B:241:LEU:HD12	2.33	0.43
1:B:323:THR:OG1	1:B:324:GLU:OE1	2.36	0.43
4:A:1308:NAG:H82	1:C:557:LYS:HD3	2.00	0.43
1:B:995:LEU:HD23	1:B:995:LEU:HA	1.87	0.43
1:C:372:ALA:HB3	1:C:374:PHE:CE1	2.54	0.43
1:C:1126:ASP:OD1	1:C:1126:ASP:N	2.51	0.43
1:A:392:PHE:HB3	1:A:395:VAL:HB	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:409:GLN:HA	1:B:414:GLN:HG2	2.00	0.43
1:A:489:PHE:CE2	1:A:491:LEU:HB2	2.53	0.43
1:A:346:ARG:HH22	1:A:448:ASN:HD21	1.67	0.42
1:A:698:LEU:HD22	1:B:872:TYR:OH	2.18	0.42
1:A:743:GLY:H	1:A:976:LEU:HD12	1.84	0.42
1:B:702:ASN:O	1:C:788:TYR:HA	2.19	0.42
1:A:353:TRP:CD1	1:A:353:TRP:H	2.37	0.42
1:A:581:LEU:HD12	1:A:581:LEU:HA	1.91	0.42
1:C:984:ASP:OD1	1:C:984:ASP:N	2.52	0.42
1:A:40:ASP:OD1	1:A:40:ASP:N	2.52	0.42
1:A:133:PHE:HB3	1:A:162:SER:O	2.19	0.42
1:A:303:LEU:HD12	1:A:308:VAL:HG12	2.00	0.42
1:A:351:TYR:HE1	1:A:452:TRP:HB2	1.85	0.42
1:B:461:LEU:HD11	1:B:467:ASP:HB3	2.00	0.42
1:B:698:LEU:HD22	1:C:872:TYR:CZ	2.54	0.42
1:A:105:ILE:HD11	1:A:239:GLN:HB2	2.01	0.42
1:B:106:PHE:HD1	1:B:238:PHE:HB2	1.85	0.42
1:C:452:TRP:CE2	1:C:493:SER:HB3	2.54	0.42
1:A:982:ARG:NH1	1:C:516:LEU:HD23	2.34	0.42
1:B:316:SER:O	1:B:594:VAL:HG12	2.20	0.42
1:B:344:ALA:HB3	1:B:347:PHE:CE1	2.55	0.42
1:C:379:CYS:HA	1:C:432:CYS:HA	2.01	0.42
1:A:534:LYS:HE2	1:A:534:LYS:HB3	1.78	0.42
1:B:698:LEU:HD11	1:C:868:MET:HG2	2.01	0.42
1:B:970:GLY:O	1:B:994:ARG:NE	2.53	0.42
1:A:666:GLY:O	1:A:669:VAL:HG22	2.19	0.42
1:B:342:PHE:CZ	1:B:510:VAL:HG11	2.54	0.42
1:C:500:TYR:HB3	1:C:504:HIS:HB2	2.02	0.42
1:C:421:TYR:CD1	1:C:457:ARG:HB3	2.54	0.42
1:C:551:LEU:HB3	1:C:584:LEU:HD13	2.02	0.42
1:A:500:TYR:HB3	1:A:504:HIS:HB2	2.01	0.42
1:B:631:THR:O	1:B:633:ARG:HD2	2.19	0.42
1:C:54:LEU:HD12	1:C:195:LYS:HD2	2.02	0.42
1:A:551:LEU:HD22	1:A:584:LEU:HD13	2.02	0.41
1:A:732:LYS:HD2	1:A:770:ALA:HB1	2.02	0.41
1:B:172:SER:O	1:B:172:SER:OG	2.31	0.41
1:C:1100:HIS:ND1	4:C:1305:NAG:H5	2.35	0.41
1:A:403:LYS:HE2	1:A:403:LYS:HB3	1.88	0.41
1:A:702:ASN:O	1:B:788:TYR:HA	2.20	0.41
1:B:435:ALA:HB2	1:B:509:VAL:HG22	2.03	0.41
1:B:758:PHE:O	1:B:762:LEU:HG	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:958:LEU:O	1:B:962:VAL:HG22	2.20	0.41
4:B:1302:NAG:H5	1:C:795:TYR:CE1	2.55	0.41
1:C:456:PHE:HB2	1:C:490:PRO:HB3	2.02	0.41
1:C:568:ILE:HG23	1:C:569:VAL:HG23	2.02	0.41
1:A:335:LEU:HB2	1:A:361:CYS:HB2	2.02	0.41
1:A:357:ARG:HH21	1:A:359:SER:HB3	1.85	0.41
1:A:358:ILE:HB	1:A:395:VAL:HG13	2.03	0.41
1:B:902:ALA:HB1	1:B:912:GLN:HB2	2.01	0.41
1:C:709:ASN:OD1	1:C:709:ASN:N	2.54	0.41
1:C:1089:PRO:HD3	1:C:1094:PHE:CE2	2.56	0.41
1:B:98:SER:HB2	1:B:100:ILE:HG12	2.02	0.41
1:B:379:CYS:HA	1:B:432:CYS:HA	2.02	0.41
1:C:402:ILE:HB	1:C:406:GLU:CD	2.46	0.41
1:C:518:HIS:HB3	1:C:564:PHE:HE2	1.86	0.41
1:A:628:LEU:HD23	1:A:628:LEU:HA	1.81	0.41
1:A:1096:SER:HB2	1:A:1101:TRP:CD2	2.56	0.41
1:B:437:ASN:OD1	1:B:505:GLN:HB3	2.21	0.41
1:B:567:ASP:OD1	1:B:567:ASP:N	2.54	0.41
1:B:709:ASN:OD1	1:B:709:ASN:N	2.54	0.41
1:C:807:ASP:HB3	1:C:810:LYS:HE3	2.02	0.41
1:C:927:ASN:O	1:C:930:ILE:HG22	2.21	0.41
1:A:400:PHE:HB2	1:A:509:VAL:HB	2.03	0.41
1:A:478:LYS:H	1:A:478:LYS:HG2	1.62	0.41
1:A:807:ASP:HB3	1:A:810:LYS:HE3	2.03	0.41
1:B:1093:VAL:HG13	1:C:899:MET:HE1	2.02	0.41
1:B:1096:SER:HB2	1:B:1101:TRP:CD2	2.56	0.41
1:A:758:PHE:O	1:A:762:LEU:HG	2.22	0.40
1:B:436:TRP:CE2	1:B:508:ARG:HB3	2.56	0.40
1:B:478:LYS:H	1:B:478:LYS:HG2	1.72	0.40
1:C:794:LYS:HD3	1:C:794:LYS:HA	1.82	0.40
1:B:329:PHE:HD1	1:B:330:PRO:HD2	1.87	0.40
1:B:885:TRP:HE3	1:B:885:TRP:H	1.65	0.40
1:C:137:ASN:HA	1:C:160:TYR:HA	2.02	0.40
1:C:200:TYR:CZ	1:C:230:PRO:HB3	2.56	0.40
1:B:574:ALA:HA	1:B:585:ASP:HA	2.04	0.40
1:C:138:ASP:HA	1:C:139:PRO:HD2	1.82	0.40
1:C:985:PRO:N	1:C:986:PRO:HD2	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	997/1248 (80%)	941 (94%)	55 (6%)	1 (0%)	48	71
1	B	980/1248 (78%)	904 (92%)	74 (8%)	2 (0%)	43	65
1	C	997/1248 (80%)	928 (93%)	65 (6%)	4 (0%)	30	54
All	All	2974/3744 (79%)	2773 (93%)	194 (6%)	7 (0%)	44	65

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	160	TYR
1	A	160	TYR
1	C	283	GLY
1	C	332	VAL
1	B	330	PRO
1	C	160	TYR
1	C	523	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	888/1087 (82%)	888 (100%)	0	100	100
1	B	876/1087 (81%)	875 (100%)	1 (0%)	88	96
1	C	888/1087 (82%)	887 (100%)	1 (0%)	88	96
All	All	2652/3261 (81%)	2650 (100%)	2 (0%)	87	96

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	330	PRO
1	C	332	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	239	GLN
1	A	314	GLN
1	A	360	ASN
1	A	448	ASN
1	A	504	HIS
1	A	563	GLN
1	A	612	GLN
1	A	689	GLN
1	A	918	ASN
1	A	927	ASN
1	A	1009	GLN
1	A	1053	GLN
1	B	121	ASN
1	B	314	GLN
1	B	334	ASN
1	B	339	HIS
1	B	417	ASN
1	B	606	GLN
1	B	612	GLN
1	B	689	GLN
1	B	754	GLN
1	B	761	GLN
1	B	786	GLN
1	B	918	ASN
1	B	948	GLN
1	B	954	ASN
1	B	1053	GLN
1	B	1087	HIS
1	B	1105	GLN
1	C	137	ASN
1	C	360	ASN
1	C	445	HIS
1	C	563	GLN
1	C	606	GLN

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Mol	Chain	Res	Type
1	C	640	ASN
1	C	702	ASN
1	C	900	GLN
1	C	918	ASN
1	C	952	ASN
1	C	954	ASN
1	C	1087	HIS

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 ⓘ

9 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	1,2	14,14,15	0.23	0	17,19,21	0.56	0
2	NAG	D	2	2	14,14,15	0.24	0	17,19,21	0.47	0
2	NAG	E	1	1,2	14,14,15	0.21	0	17,19,21	0.45	0
2	NAG	E	2	2	14,14,15	0.27	0	17,19,21	0.50	0
2	NAG	F	1	1,2	14,14,15	0.40	0	17,19,21	0.76	0
2	NAG	F	2	2	14,14,15	0.36	0	17,19,21	0.98	1 (5%)
3	NAG	G	1	1,3	14,14,15	0.38	0	17,19,21	1.29	2 (11%)
3	NAG	G	2	3	14,14,15	0.39	0	17,19,21	0.56	0
3	BMA	G	3	3	11,11,12	0.29	0	15,15,17	0.54	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	1,2	-	4/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1
3	NAG	G	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	BMA	G	3	3	-	1/2/19/22	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(°)
3	G	1	NAG	C1-C2-N2	3.28	115.60	110.43
2	F	2	NAG	C2-N2-C7	3.16	127.14	122.90
3	G	1	NAG	C2-N2-C7	2.85	126.72	122.90

There are no chirality outliers.

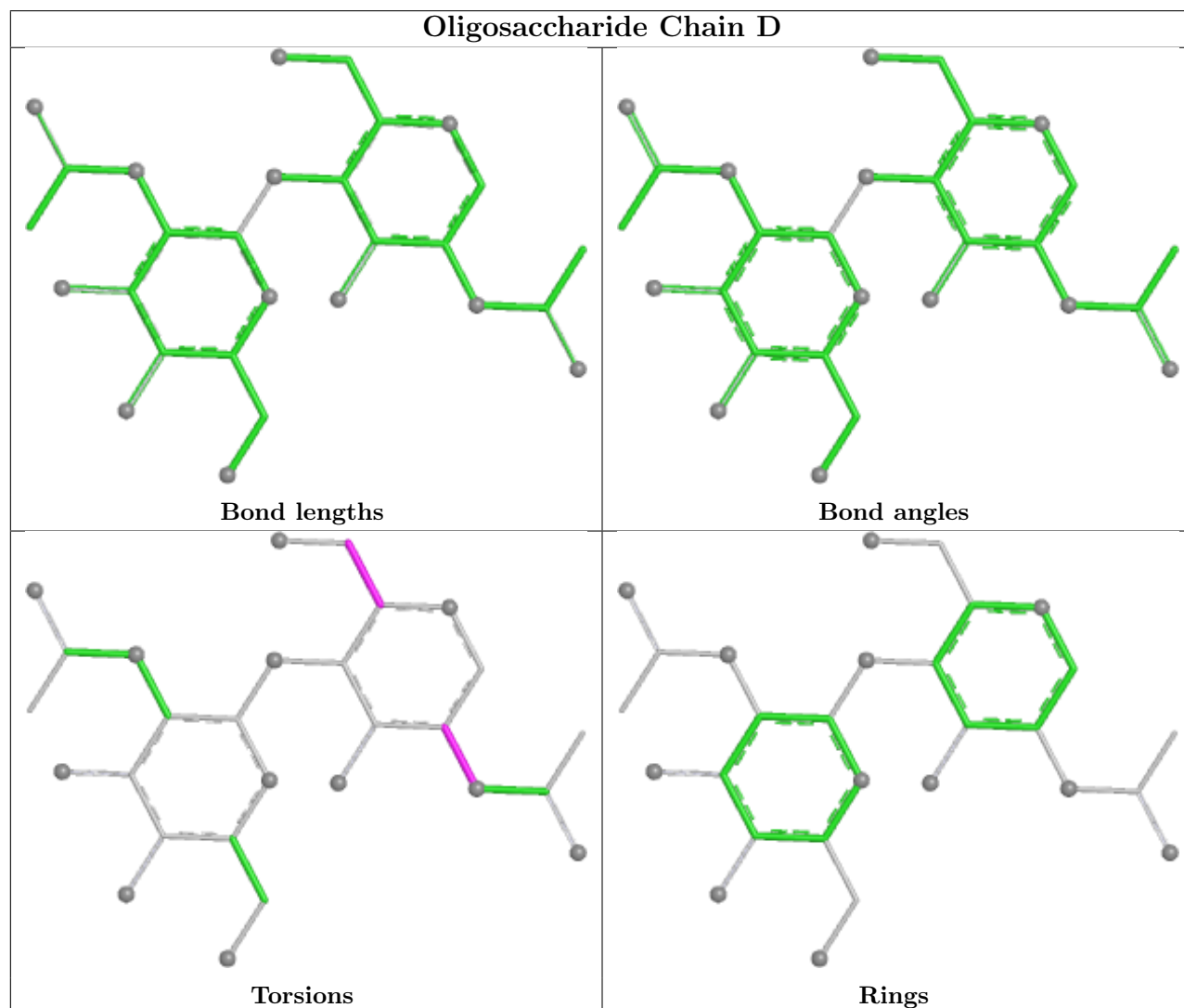
All (14) torsion outliers are listed below:

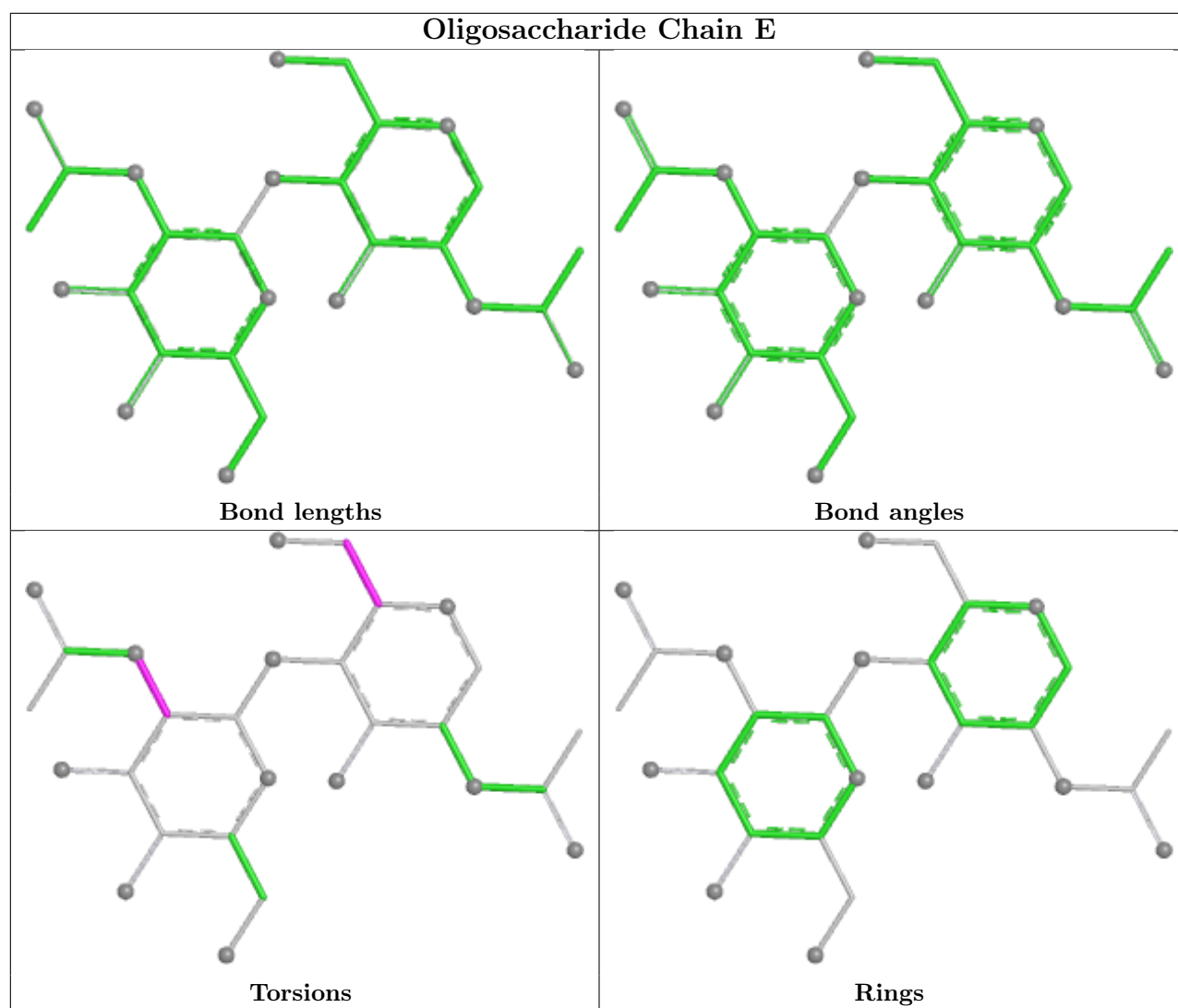
Mol	Chain	Res	Type	Atoms
3	G	1	NAG	C8-C7-N2-C2
3	G	1	NAG	O7-C7-N2-C2
3	G	3	BMA	O5-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
3	G	1	NAG	C3-C2-N2-C7
2	D	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C1-C2-N2-C7
3	G	1	NAG	C1-C2-N2-C7
2	D	1	NAG	C3-C2-N2-C7
2	E	2	NAG	C3-C2-N2-C7
2	F	2	NAG	C3-C2-N2-C7
2	D	1	NAG	O5-C5-C6-O6
2	D	1	NAG	C1-C2-N2-C7

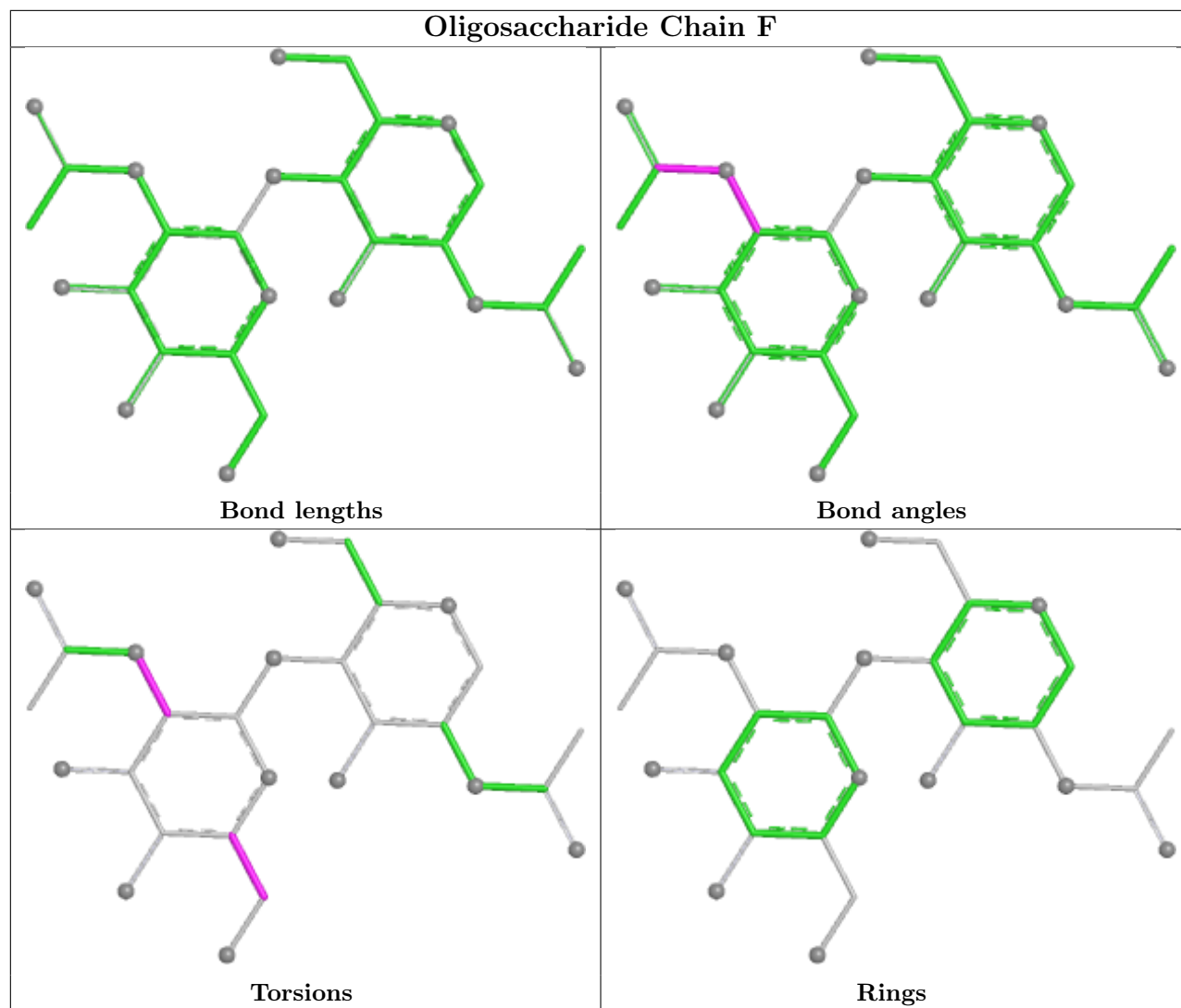
There are no ring outliers.

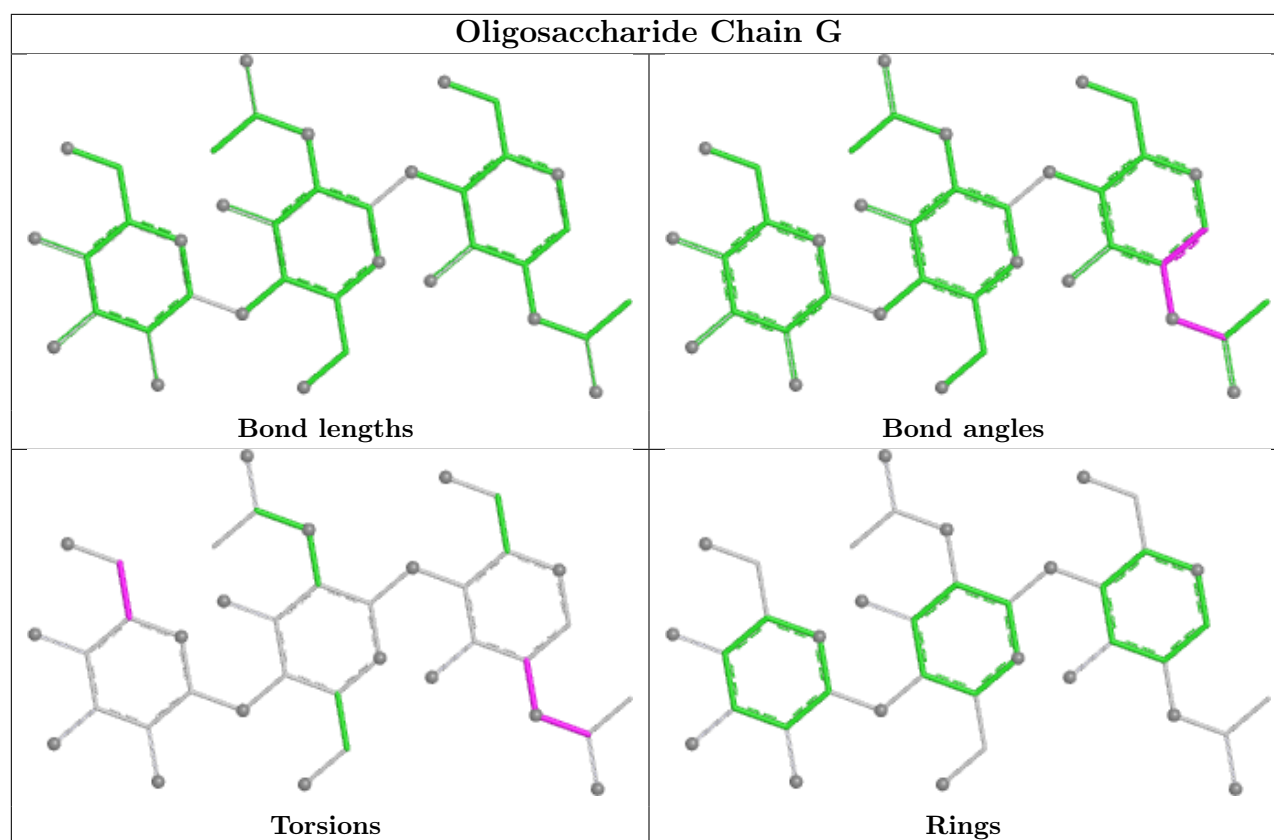
No monomer is involved in short contacts.

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](#)

26 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	B	1303	1	14,14,15	0.27	0	17,19,21	0.54	0
4	NAG	C	1305	1	14,14,15	0.22	0	17,19,21	0.53	0
4	NAG	B	1302	1	14,14,15	0.18	0	17,19,21	0.47	0
4	NAG	B	1304	1	14,14,15	0.34	0	17,19,21	0.42	0
4	NAG	A	1310	1	14,14,15	0.38	0	17,19,21	0.94	2 (11%)
4	NAG	A	1304	1	14,14,15	0.38	0	17,19,21	0.64	0
4	NAG	A	1306	1	14,14,15	0.21	0	17,19,21	0.44	0
4	NAG	B	1308	1	14,14,15	0.41	0	17,19,21	0.82	1 (5%)
4	NAG	B	1301	1	14,14,15	0.23	0	17,19,21	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	C	1304	1	14,14,15	0.32	0	17,19,21	0.40	0
4	NAG	A	1302	1	14,14,15	0.22	0	17,19,21	0.42	0
4	NAG	A	1303	1	14,14,15	0.20	0	17,19,21	0.63	0
4	NAG	C	1308	1	14,14,15	0.39	0	17,19,21	1.19	1 (5%)
4	NAG	B	1307	1	14,14,15	0.24	0	17,19,21	0.42	0
4	NAG	B	1305	1	14,14,15	0.21	0	17,19,21	0.46	0
4	NAG	C	1301	1	14,14,15	0.21	0	17,19,21	0.43	0
4	NAG	B	1306	1	14,14,15	0.20	0	17,19,21	0.44	0
4	NAG	C	1306	1	14,14,15	0.29	0	17,19,21	0.52	0
4	NAG	C	1307	1	14,14,15	0.30	0	17,19,21	0.46	0
4	NAG	A	1309	1	14,14,15	0.40	0	17,19,21	1.09	2 (11%)
4	NAG	C	1303	1	14,14,15	0.21	0	17,19,21	0.43	0
4	NAG	A	1305	1	14,14,15	0.23	0	17,19,21	0.54	0
4	NAG	A	1308	1	14,14,15	0.40	0	17,19,21	0.87	1 (5%)
4	NAG	A	1307	1	14,14,15	0.39	0	17,19,21	0.48	0
4	NAG	A	1301	1	14,14,15	0.18	0	17,19,21	0.44	0
4	NAG	C	1302	1	14,14,15	0.22	0	17,19,21	0.43	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	B	1303	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1305	1	-	3/6/23/26	0/1/1/1
4	NAG	B	1302	1	-	4/6/23/26	0/1/1/1
4	NAG	B	1304	1	-	4/6/23/26	0/1/1/1
4	NAG	A	1310	1	-	5/6/23/26	0/1/1/1
4	NAG	A	1304	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1306	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1308	1	-	1/6/23/26	0/1/1/1
4	NAG	B	1301	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1304	1	-	4/6/23/26	0/1/1/1
4	NAG	A	1302	1	-	4/6/23/26	0/1/1/1
4	NAG	A	1303	1	-	4/6/23/26	0/1/1/1
4	NAG	C	1308	1	-	1/6/23/26	0/1/1/1
4	NAG	B	1307	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1305	1	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	C	1301	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1306	1	-	1/6/23/26	0/1/1/1
4	NAG	C	1306	1	-	0/6/23/26	0/1/1/1
4	NAG	C	1307	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1309	1	-	3/6/23/26	0/1/1/1
4	NAG	C	1303	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1305	1	-	4/6/23/26	0/1/1/1
4	NAG	A	1308	1	-	3/6/23/26	0/1/1/1
4	NAG	A	1307	1	-	4/6/23/26	0/1/1/1
4	NAG	A	1301	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1302	1	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1308	NAG	C2-N2-C7	4.11	128.41	122.90
4	A	1308	NAG	C1-C2-N2	3.15	115.40	110.43
4	A	1309	NAG	C1-C2-N2	3.13	115.36	110.43
4	B	1308	NAG	C1-C2-N2	2.84	114.90	110.43
4	A	1310	NAG	C1-C2-N2	2.69	114.67	110.43
4	A	1309	NAG	C2-N2-C7	2.59	126.37	122.90
4	A	1310	NAG	C2-N2-C7	2.23	125.89	122.90

There are no chirality outliers.

All (63) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1309	NAG	C3-C2-N2-C7
4	A	1310	NAG	C1-C2-N2-C7
4	A	1310	NAG	C8-C7-N2-C2
4	A	1310	NAG	O7-C7-N2-C2
4	B	1304	NAG	C4-C5-C6-O6
4	C	1301	NAG	O5-C5-C6-O6
4	A	1305	NAG	C4-C5-C6-O6
4	C	1304	NAG	O5-C5-C6-O6
4	B	1303	NAG	O5-C5-C6-O6
4	A	1303	NAG	C4-C5-C6-O6
4	B	1301	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	A	1303	NAG	O5-C5-C6-O6
4	B	1304	NAG	O5-C5-C6-O6
4	C	1301	NAG	C4-C5-C6-O6
4	C	1304	NAG	C4-C5-C6-O6
4	A	1301	NAG	O5-C5-C6-O6
4	A	1302	NAG	C4-C5-C6-O6
4	A	1305	NAG	O5-C5-C6-O6
4	B	1302	NAG	C4-C5-C6-O6
4	A	1306	NAG	O5-C5-C6-O6
4	B	1302	NAG	O5-C5-C6-O6
4	A	1302	NAG	C8-C7-N2-C2
4	A	1302	NAG	O7-C7-N2-C2
4	A	1307	NAG	C8-C7-N2-C2
4	A	1307	NAG	O7-C7-N2-C2
4	B	1302	NAG	C8-C7-N2-C2
4	B	1302	NAG	O7-C7-N2-C2
4	B	1304	NAG	C8-C7-N2-C2
4	B	1304	NAG	O7-C7-N2-C2
4	C	1302	NAG	C8-C7-N2-C2
4	C	1302	NAG	O7-C7-N2-C2
4	C	1304	NAG	C8-C7-N2-C2
4	C	1304	NAG	O7-C7-N2-C2
4	B	1306	NAG	O5-C5-C6-O6
4	A	1302	NAG	O5-C5-C6-O6
4	A	1307	NAG	O5-C5-C6-O6
4	C	1303	NAG	C4-C5-C6-O6
4	B	1307	NAG	O5-C5-C6-O6
4	A	1306	NAG	C4-C5-C6-O6
4	B	1303	NAG	C4-C5-C6-O6
4	C	1302	NAG	O5-C5-C6-O6
4	C	1303	NAG	O5-C5-C6-O6
4	A	1301	NAG	C4-C5-C6-O6
4	A	1310	NAG	O5-C5-C6-O6
4	B	1308	NAG	O5-C5-C6-O6
4	A	1309	NAG	O5-C5-C6-O6
4	C	1305	NAG	O5-C5-C6-O6
4	B	1301	NAG	C4-C5-C6-O6
4	A	1308	NAG	C3-C2-N2-C7
4	A	1310	NAG	C3-C2-N2-C7
4	B	1305	NAG	O5-C5-C6-O6
4	A	1303	NAG	C1-C2-N2-C7
4	A	1308	NAG	C1-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
4	A	1309	NAG	C1-C2-N2-C7
4	A	1307	NAG	C4-C5-C6-O6
4	A	1303	NAG	C3-C2-N2-C7
4	A	1305	NAG	C3-C2-N2-C7
4	C	1305	NAG	C3-C2-N2-C7
4	C	1308	NAG	C3-C2-N2-C7
4	A	1305	NAG	C1-C2-N2-C7
4	C	1305	NAG	C1-C2-N2-C7
4	B	1307	NAG	C4-C5-C6-O6
4	A	1308	NAG	O7-C7-N2-C2

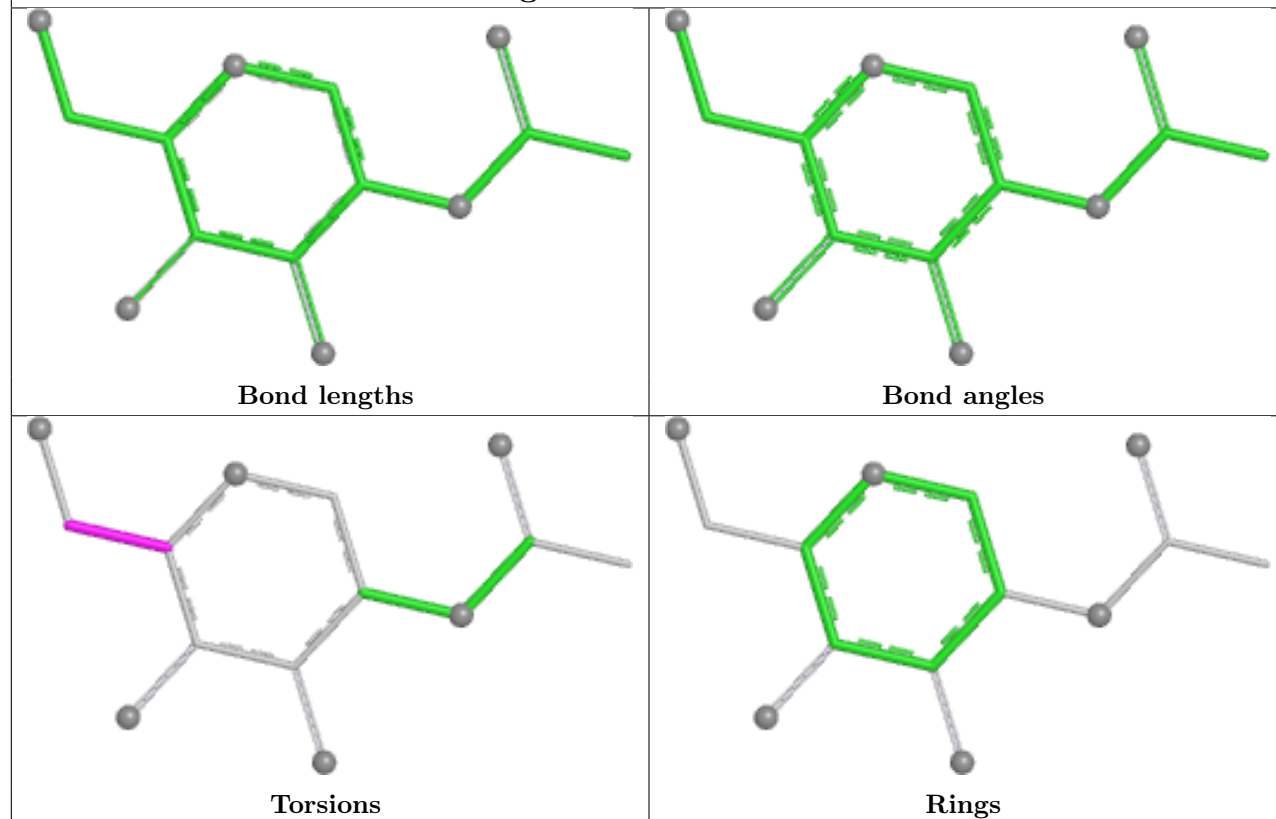
There are no ring outliers.

6 monomers are involved in 6 short contacts:

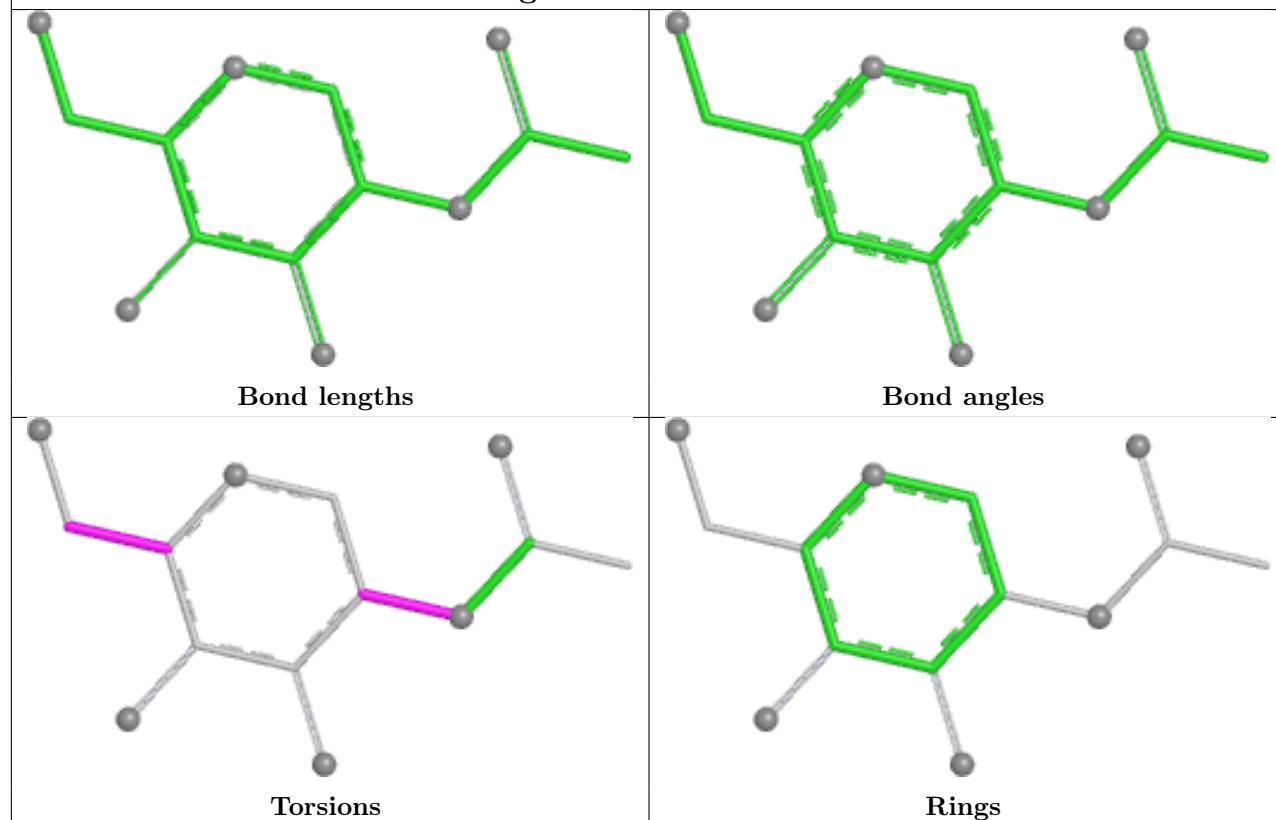
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1305	NAG	1	0
4	B	1302	NAG	1	0
4	A	1310	NAG	1	0
4	B	1308	NAG	1	0
4	A	1305	NAG	1	0
4	A	1308	NAG	1	0

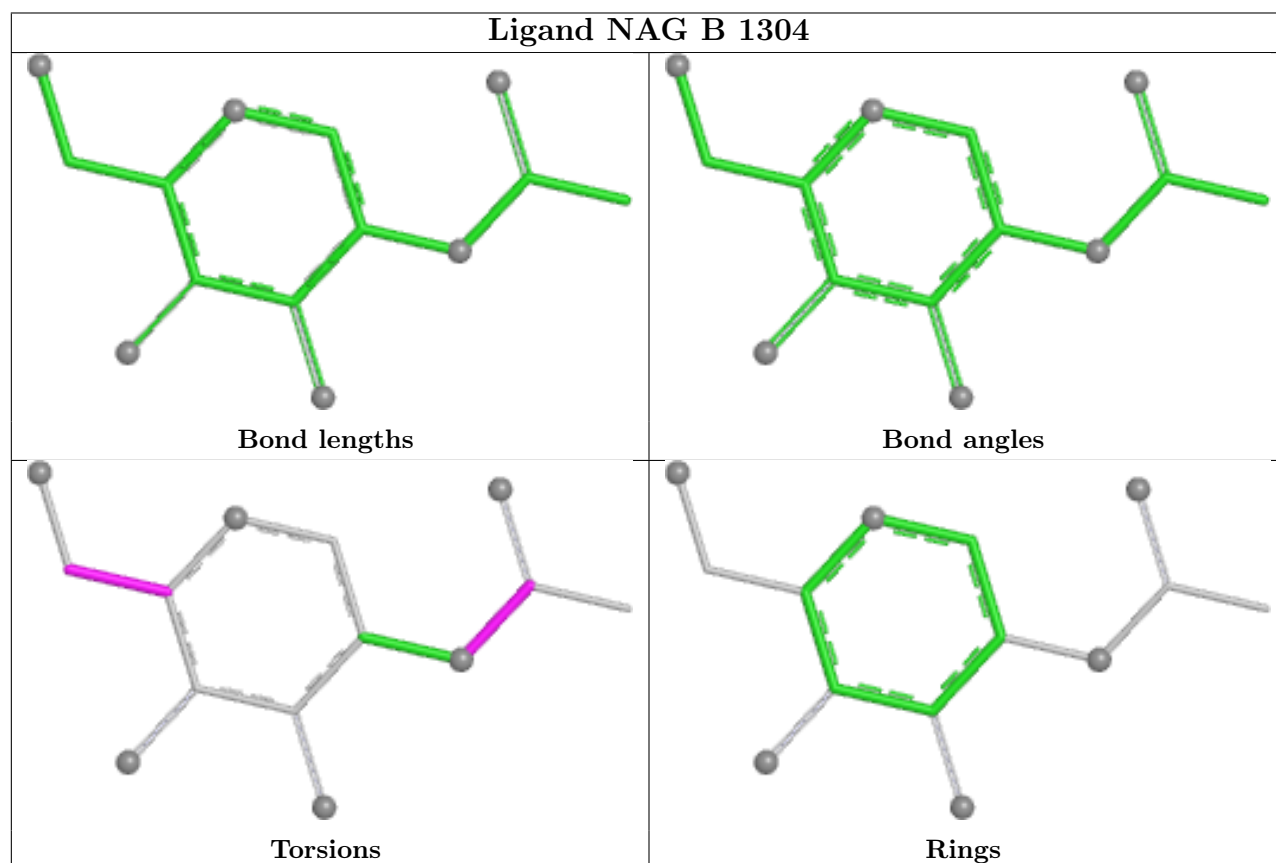
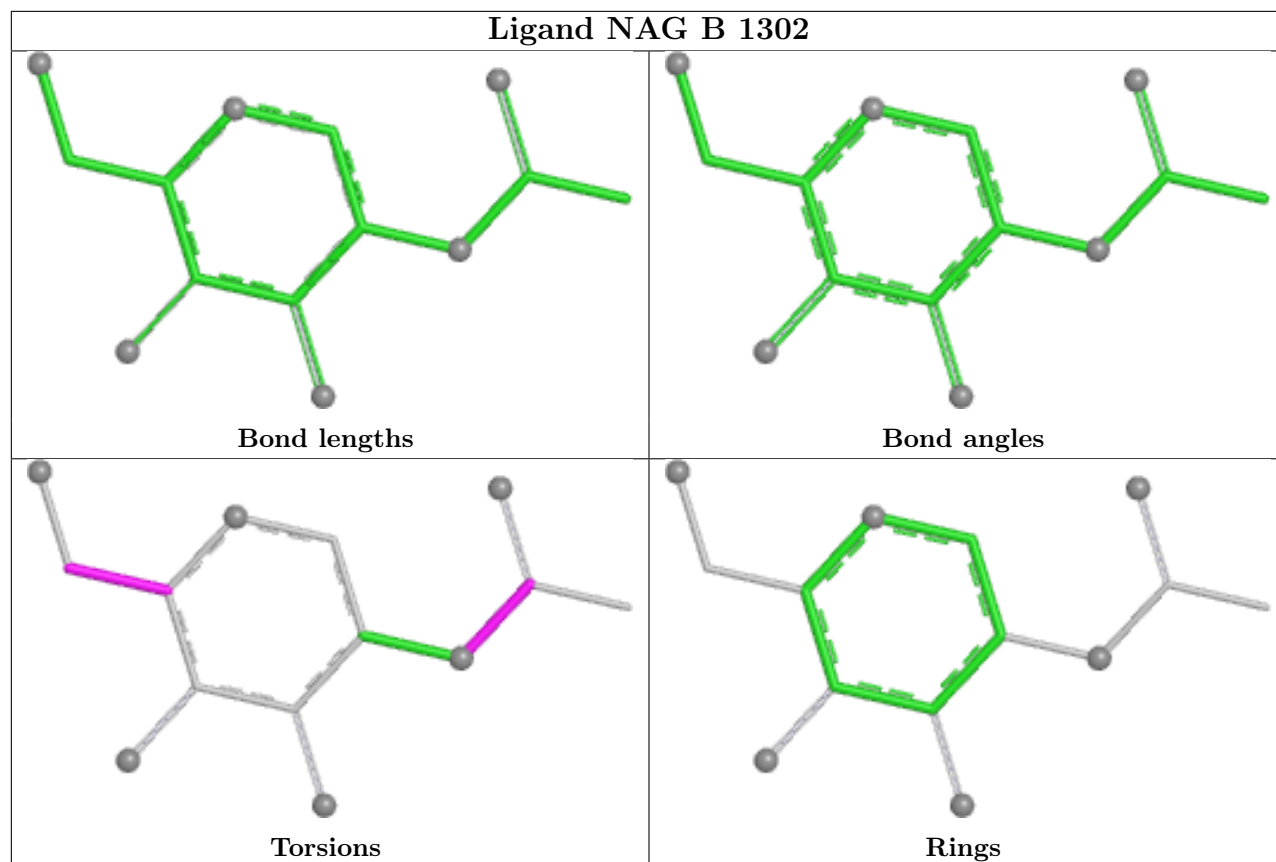
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

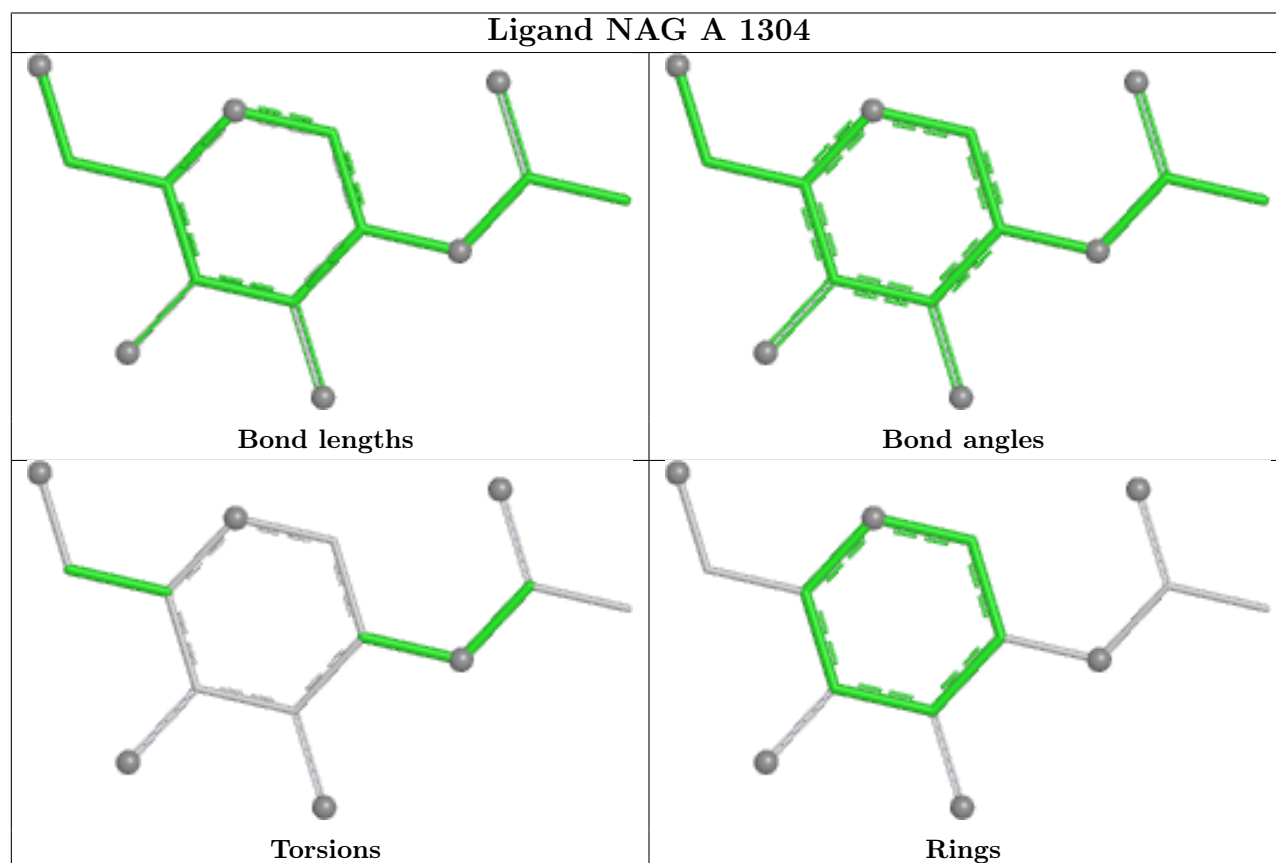
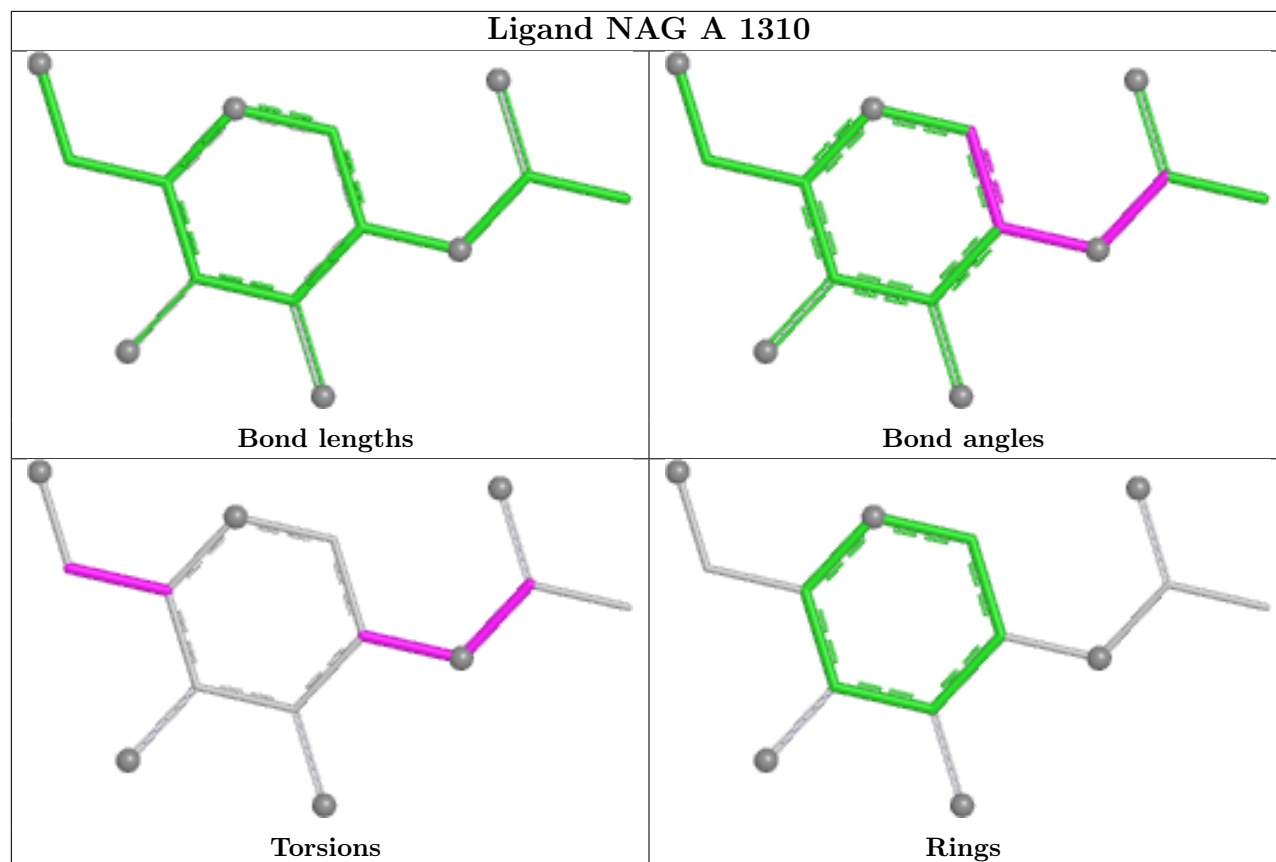
Ligand NAG B 1303



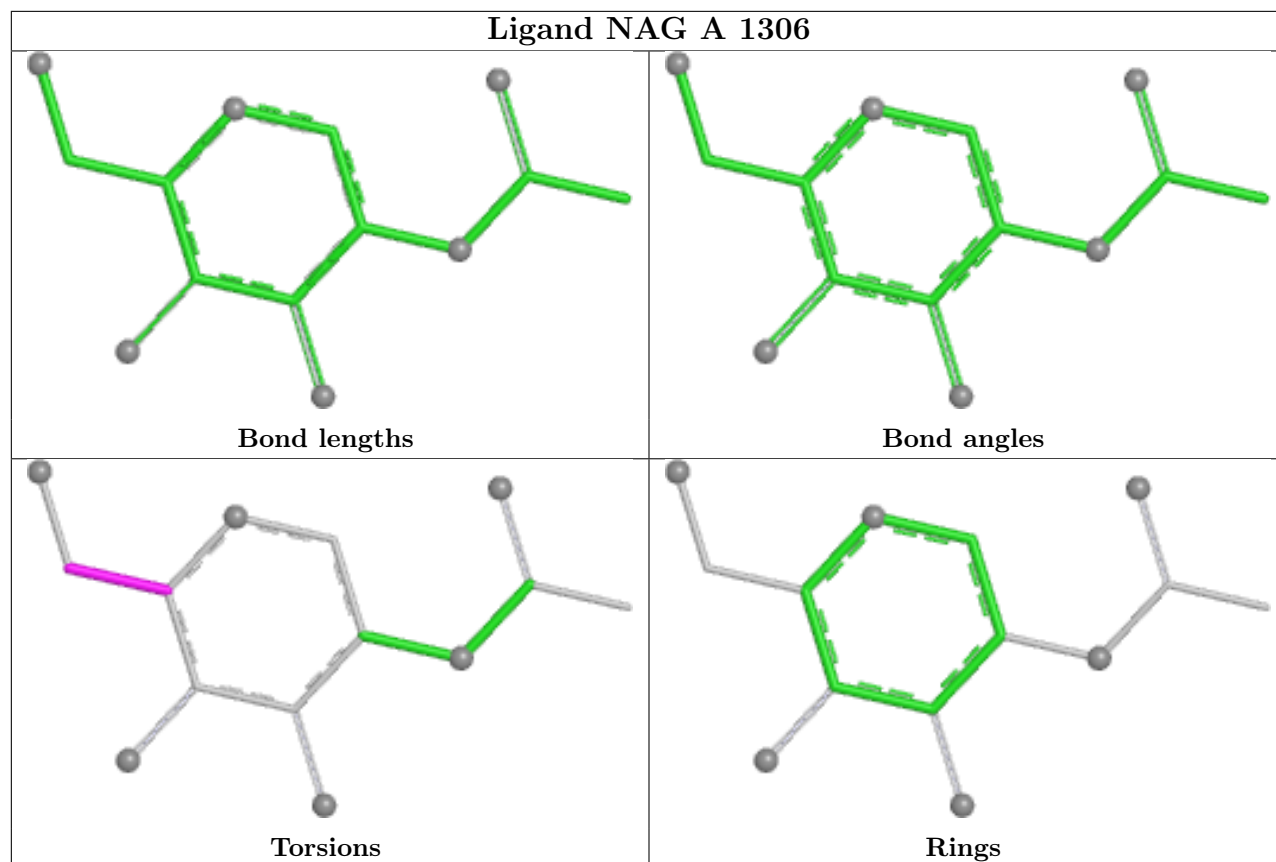
Ligand NAG C 1305



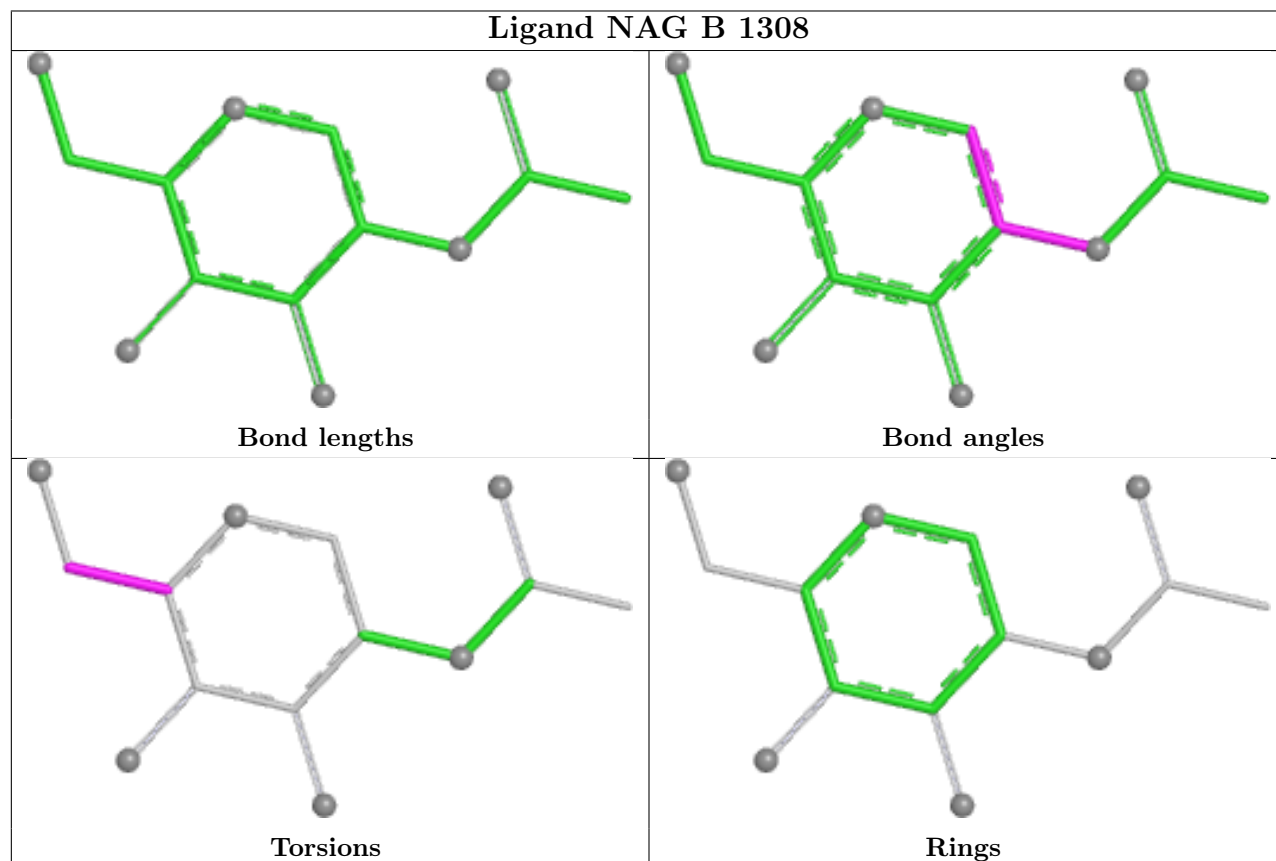




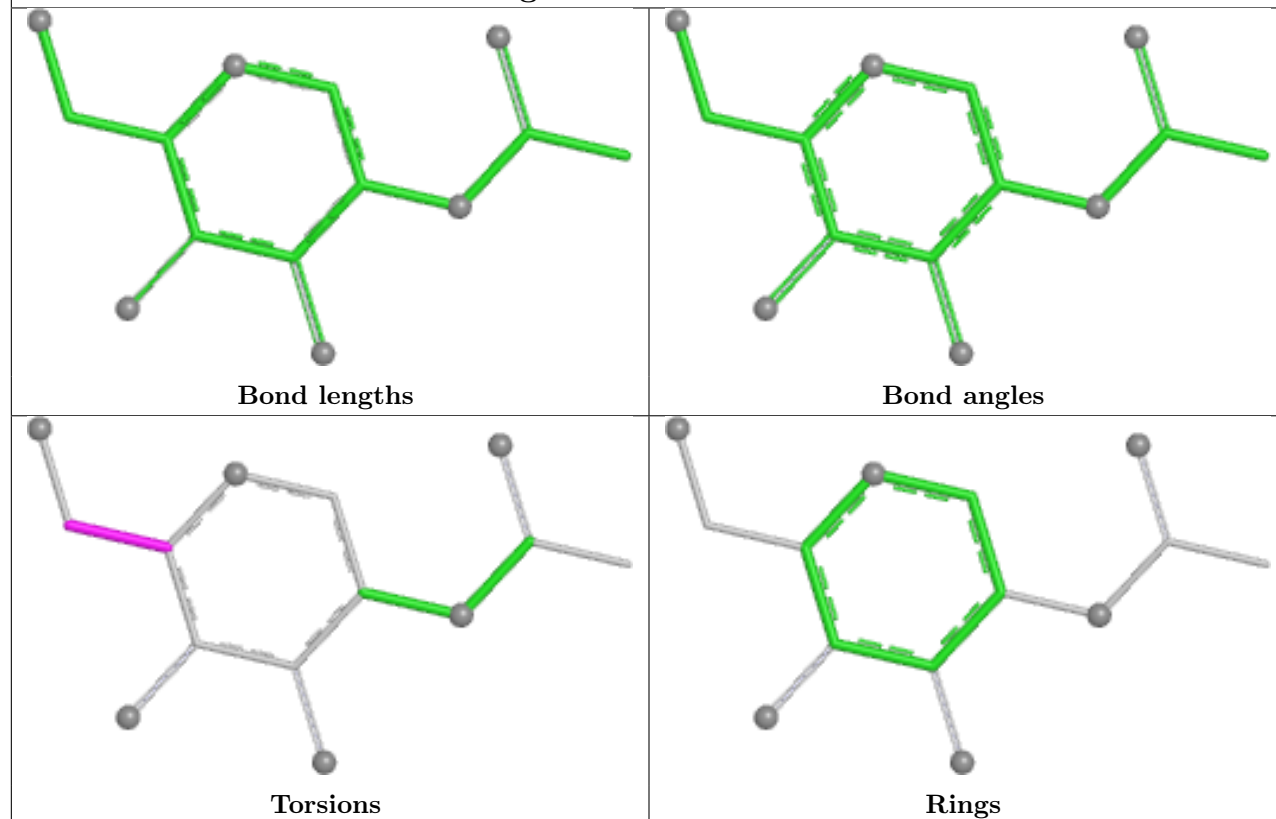
Ligand NAG A 1306



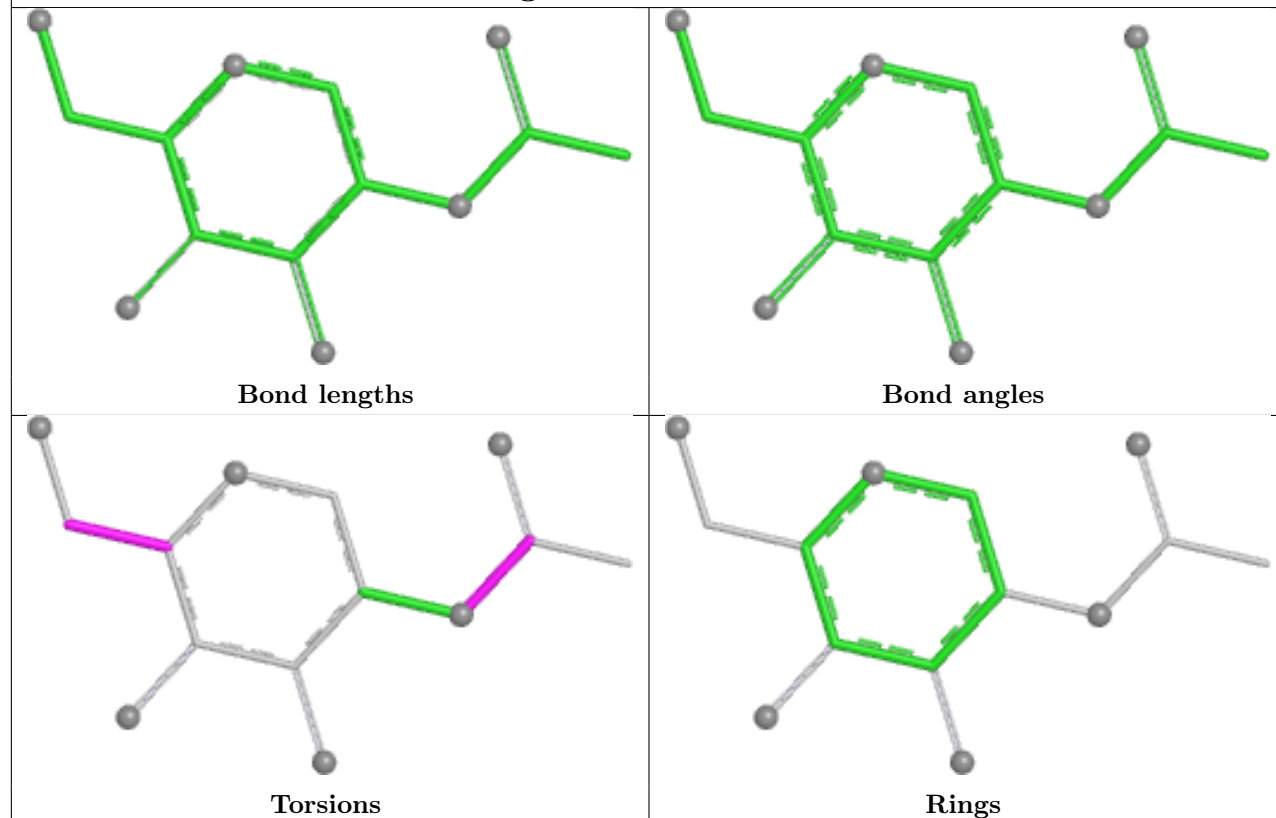
Ligand NAG B 1308



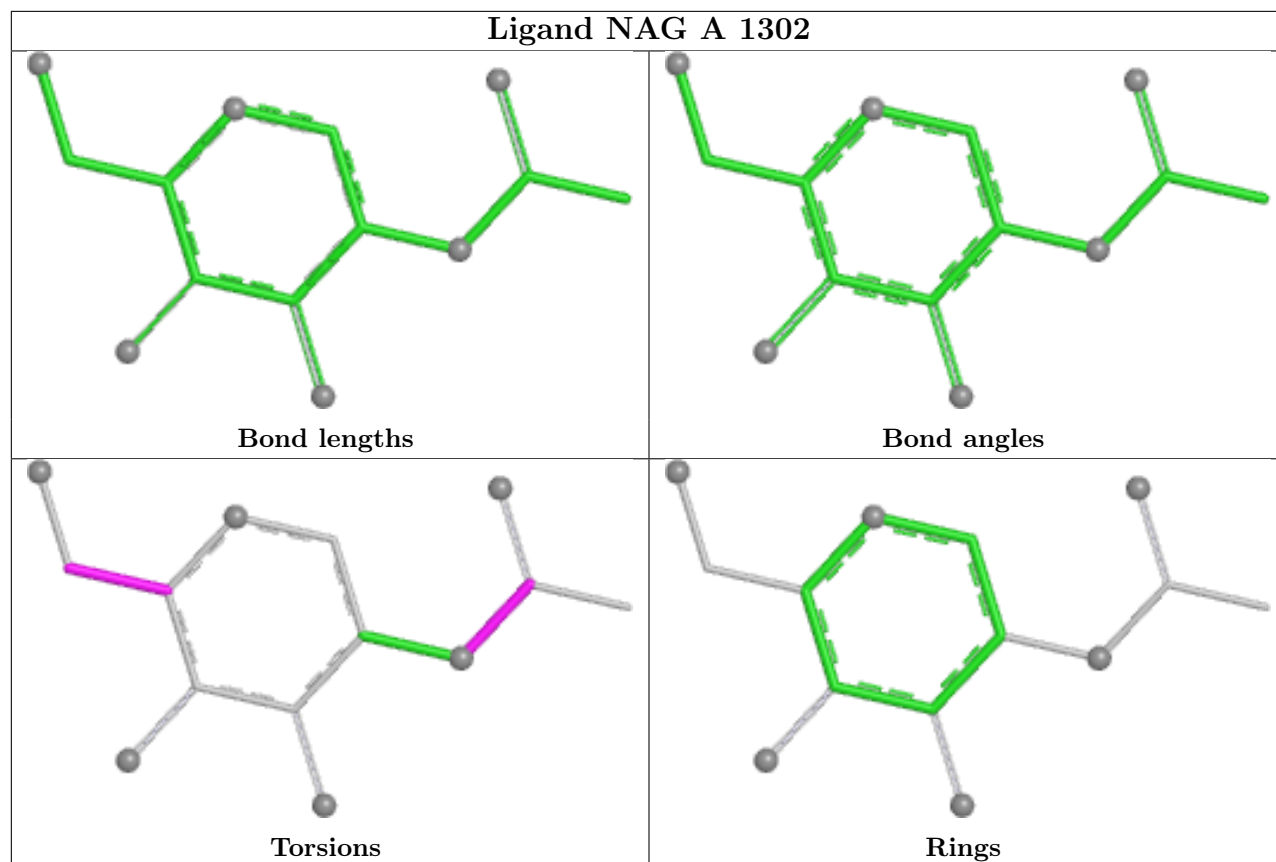
Ligand NAG B 1301



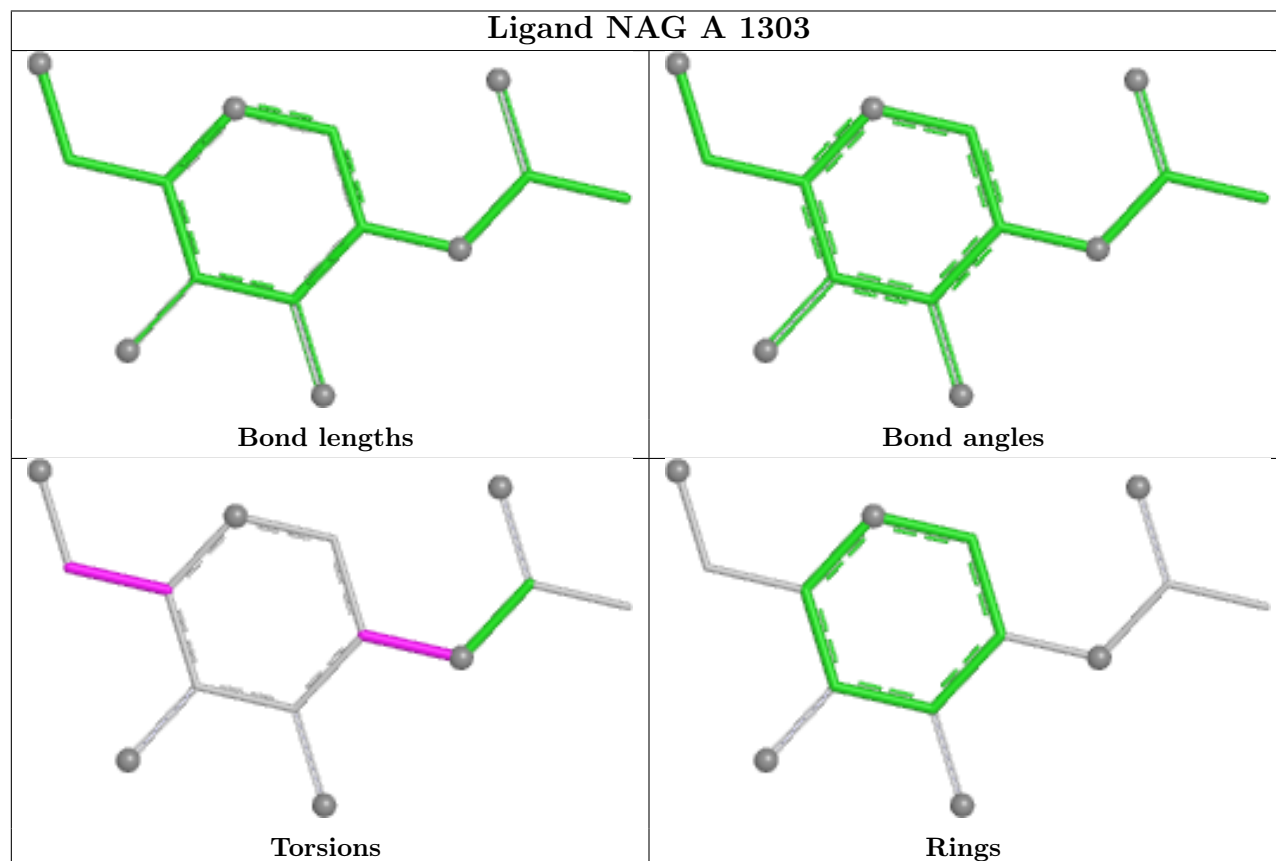
Ligand NAG C 1304

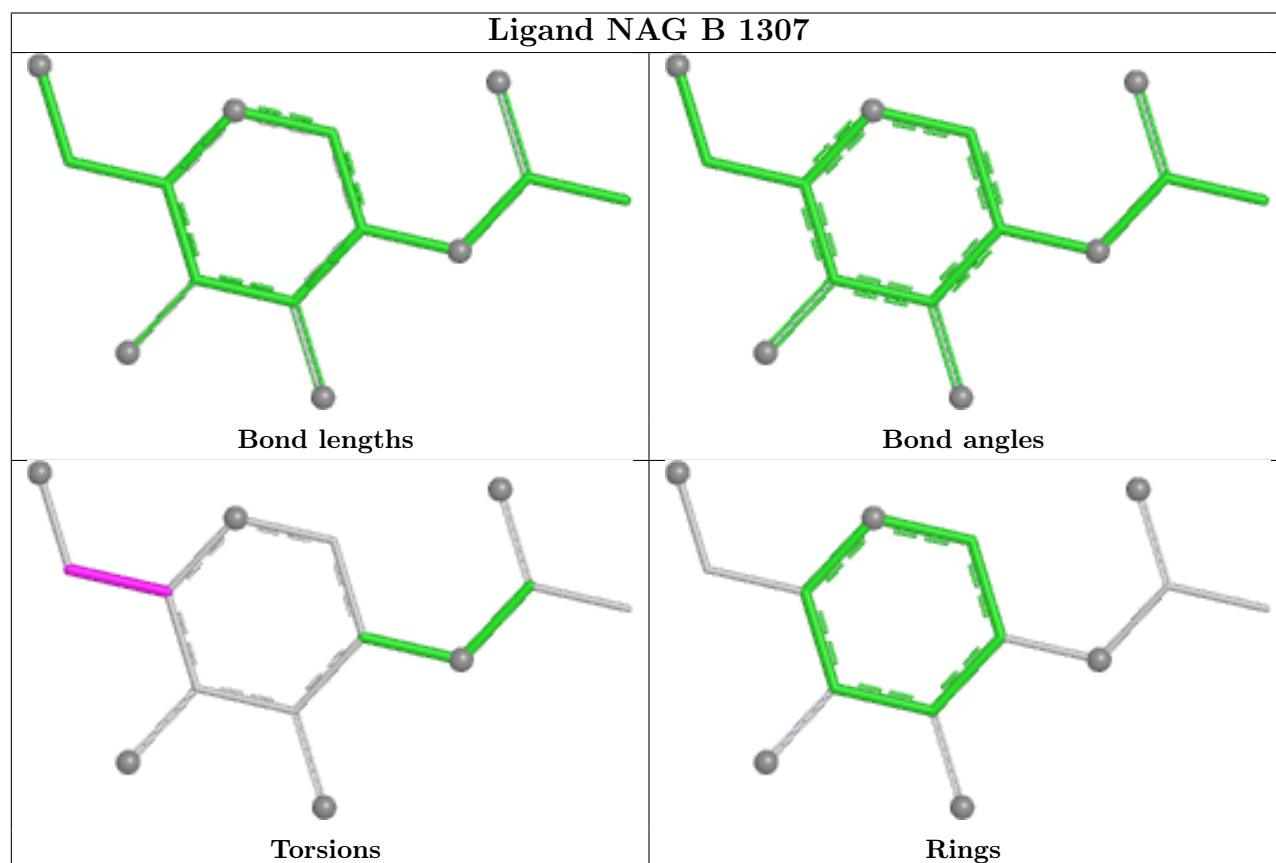
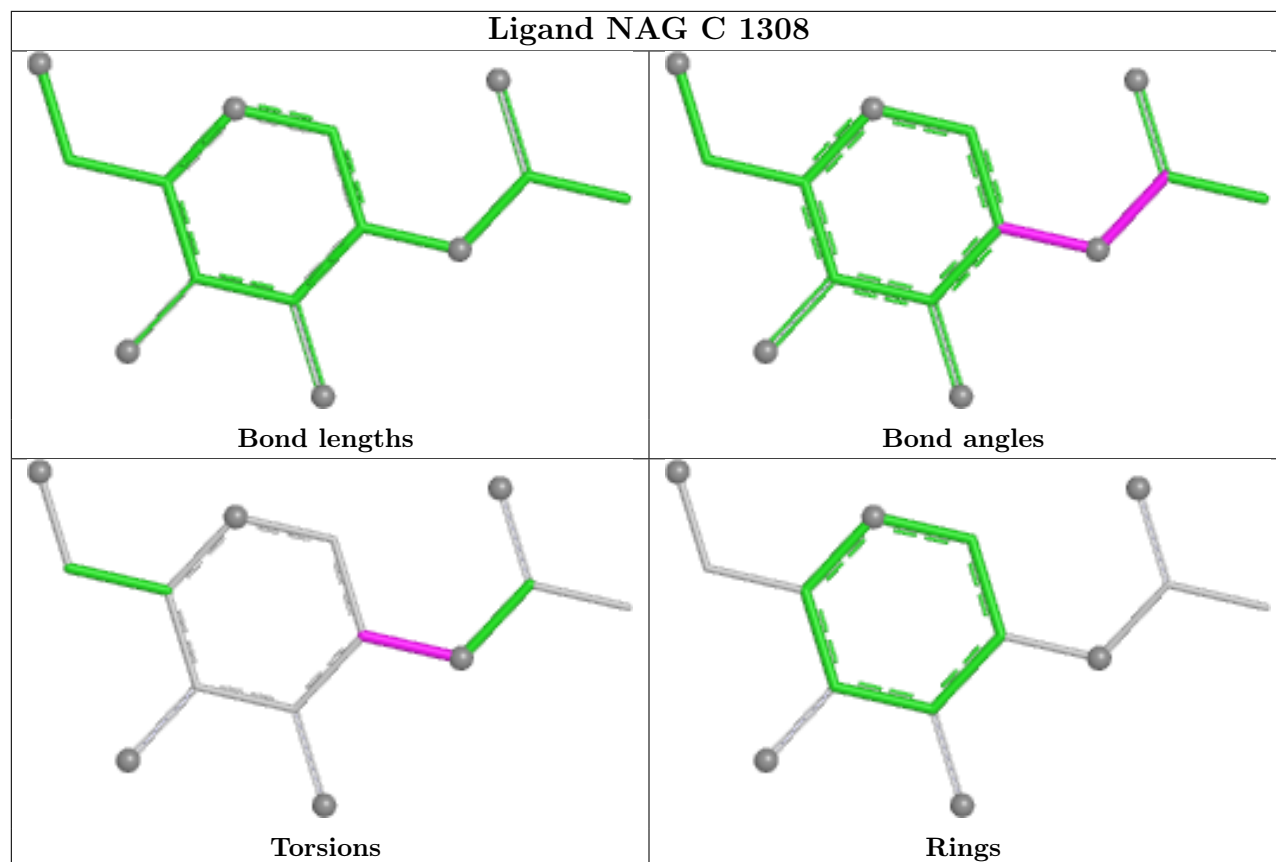


Ligand NAG A 1302

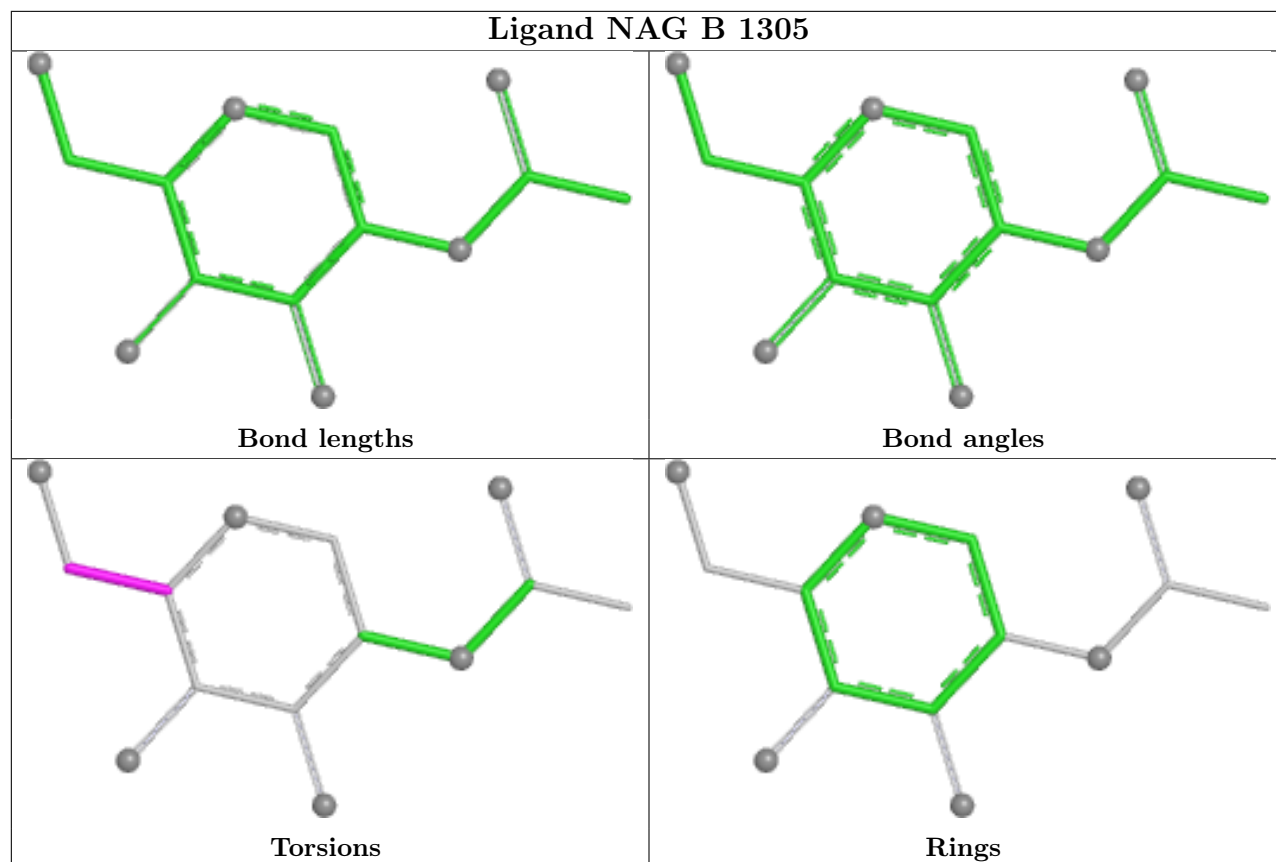


Ligand NAG A 1303

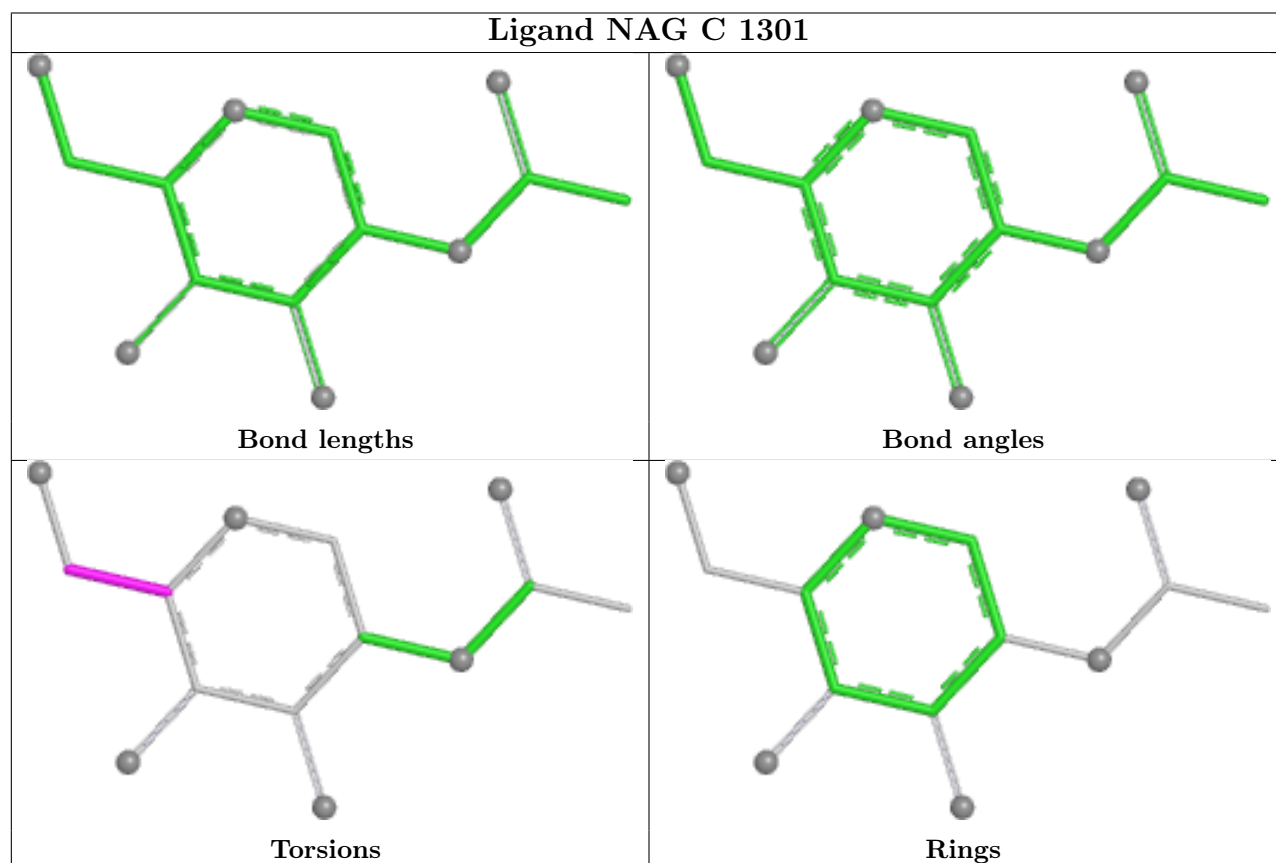




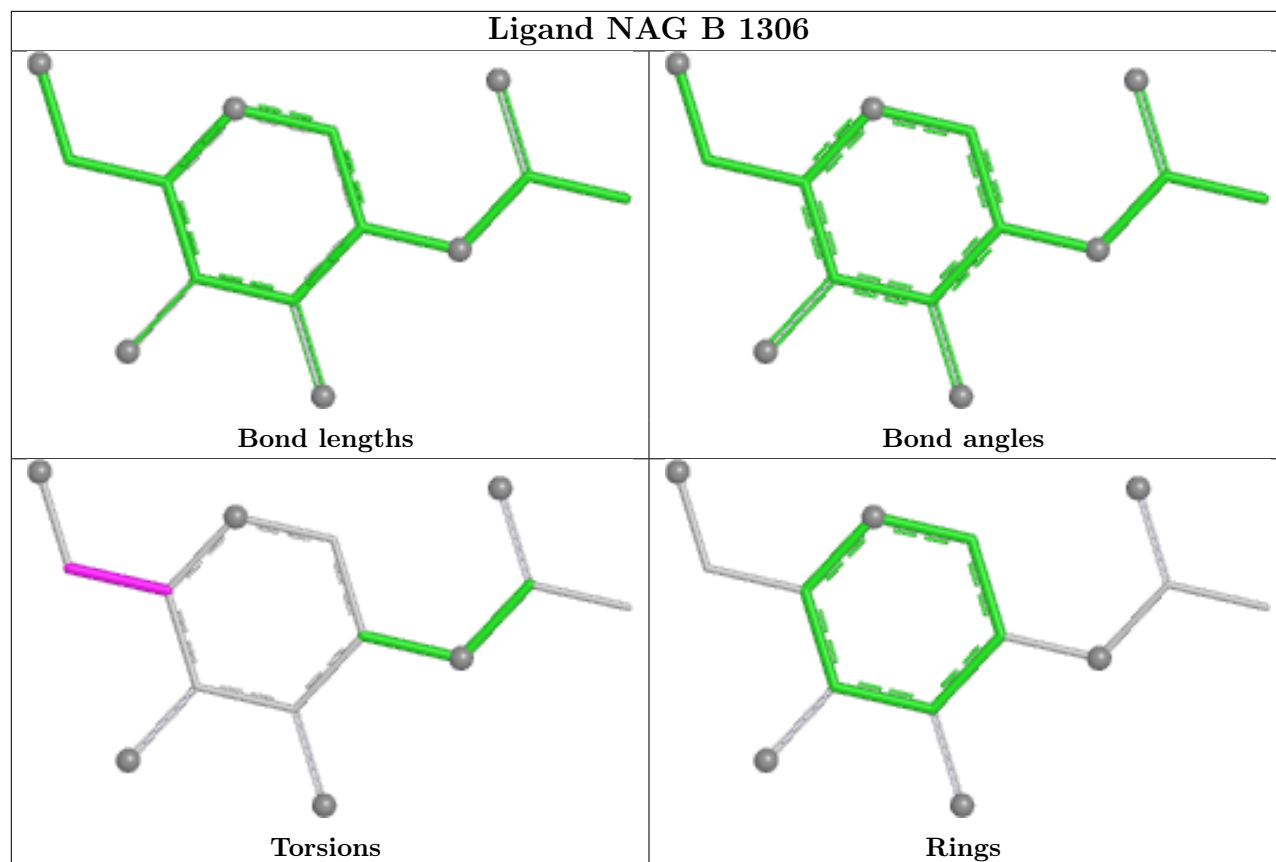
Ligand NAG B 1305



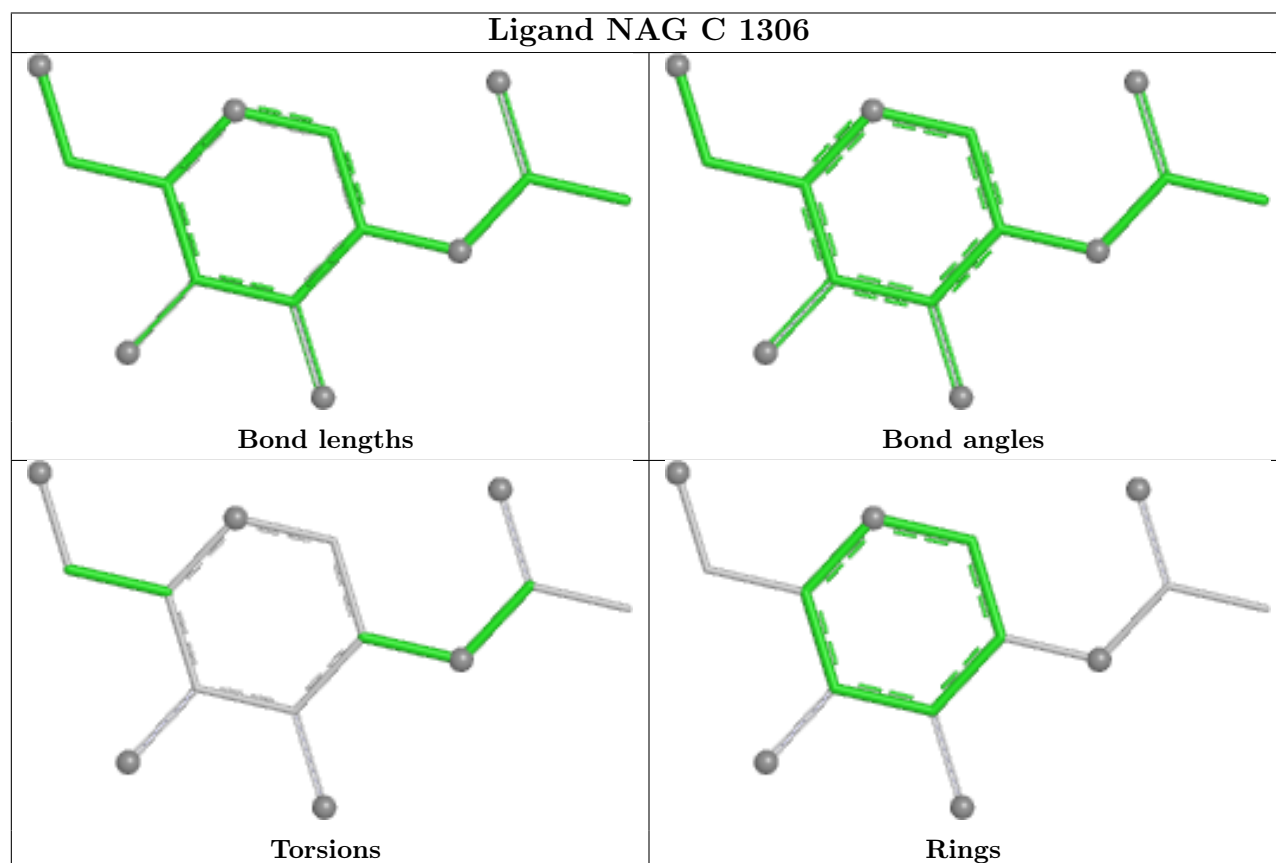
Ligand NAG C 1301



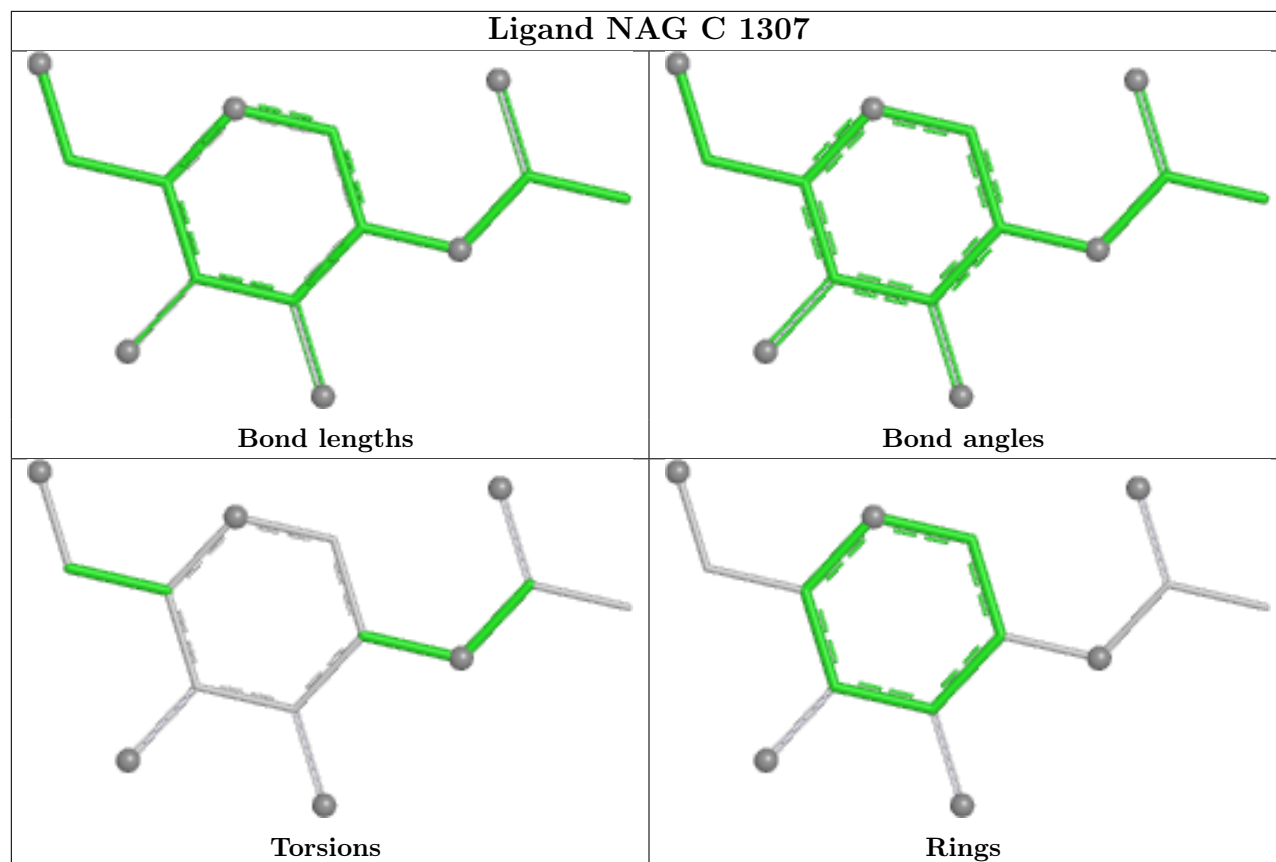
Ligand NAG B 1306



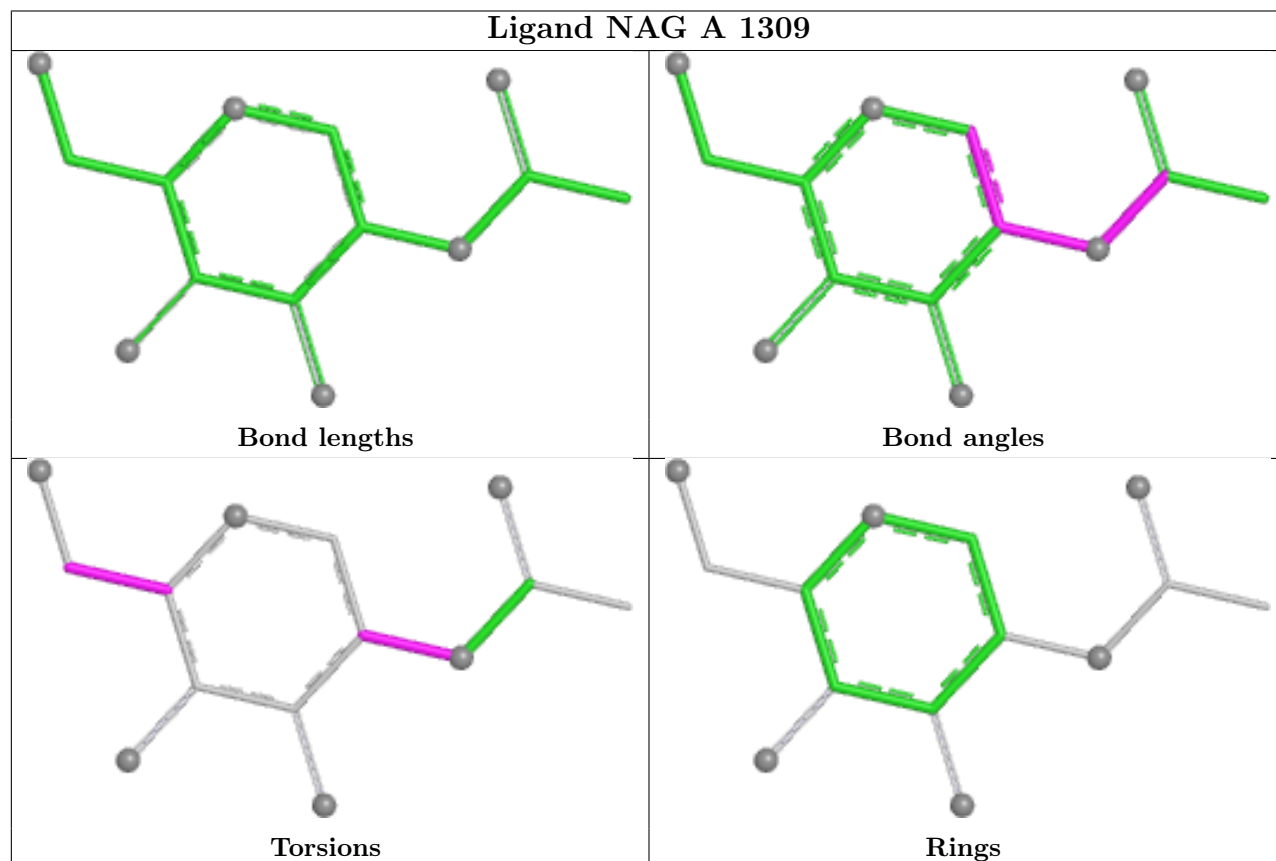
Ligand NAG C 1306



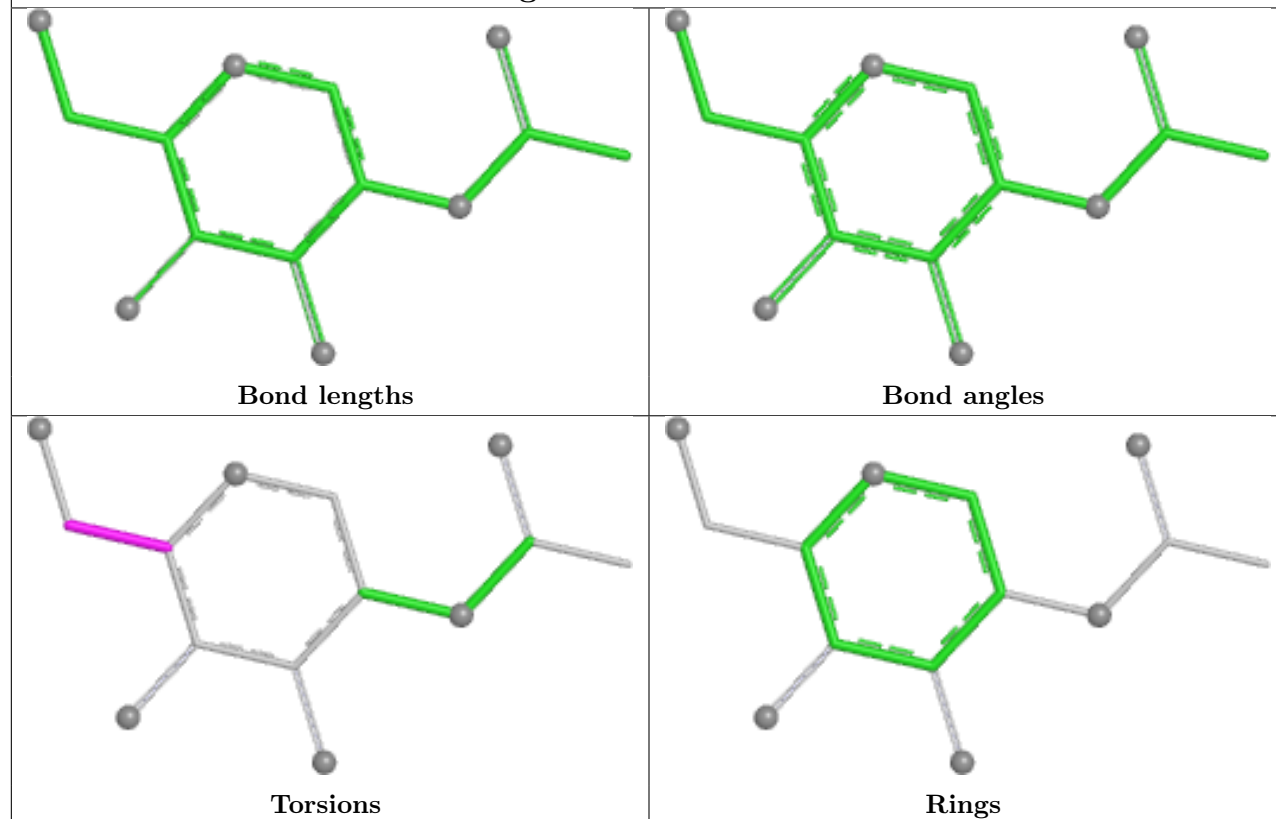
Ligand NAG C 1307



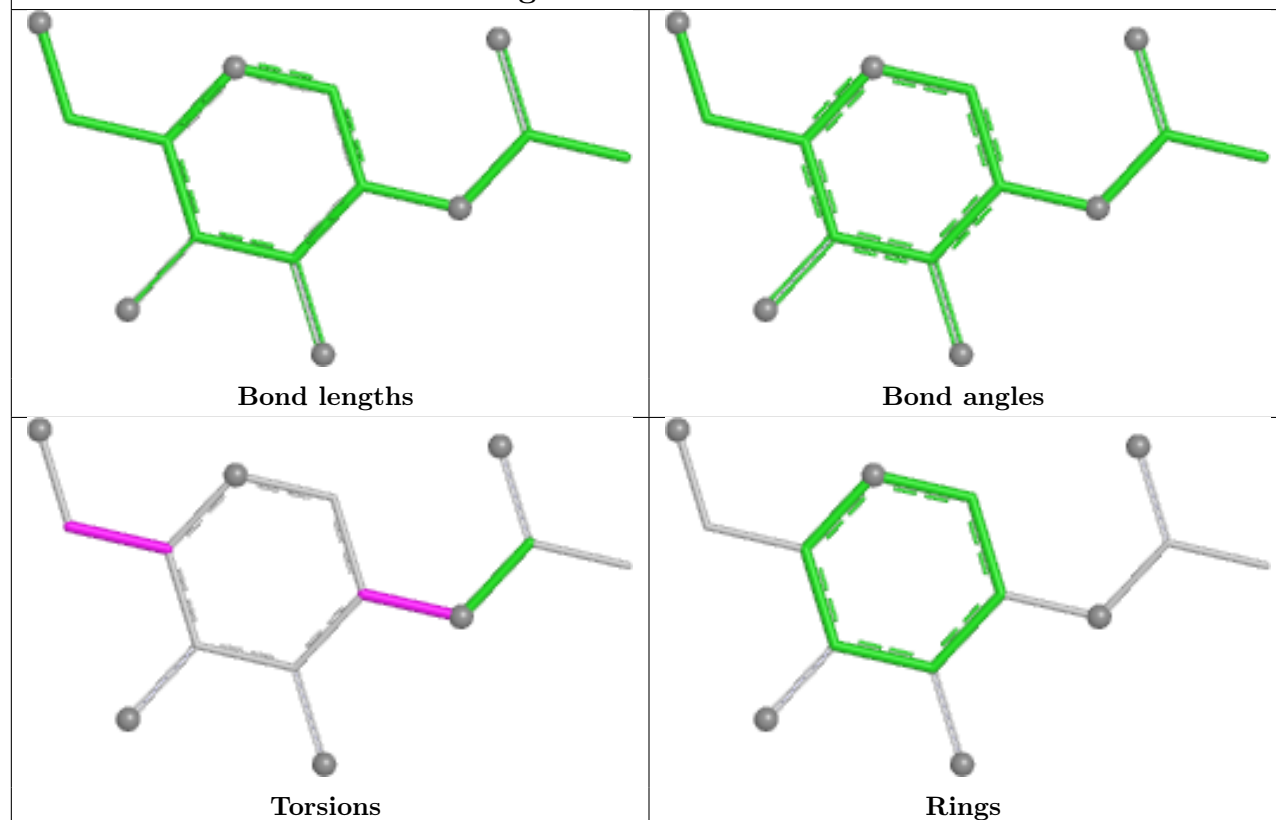
Ligand NAG A 1309

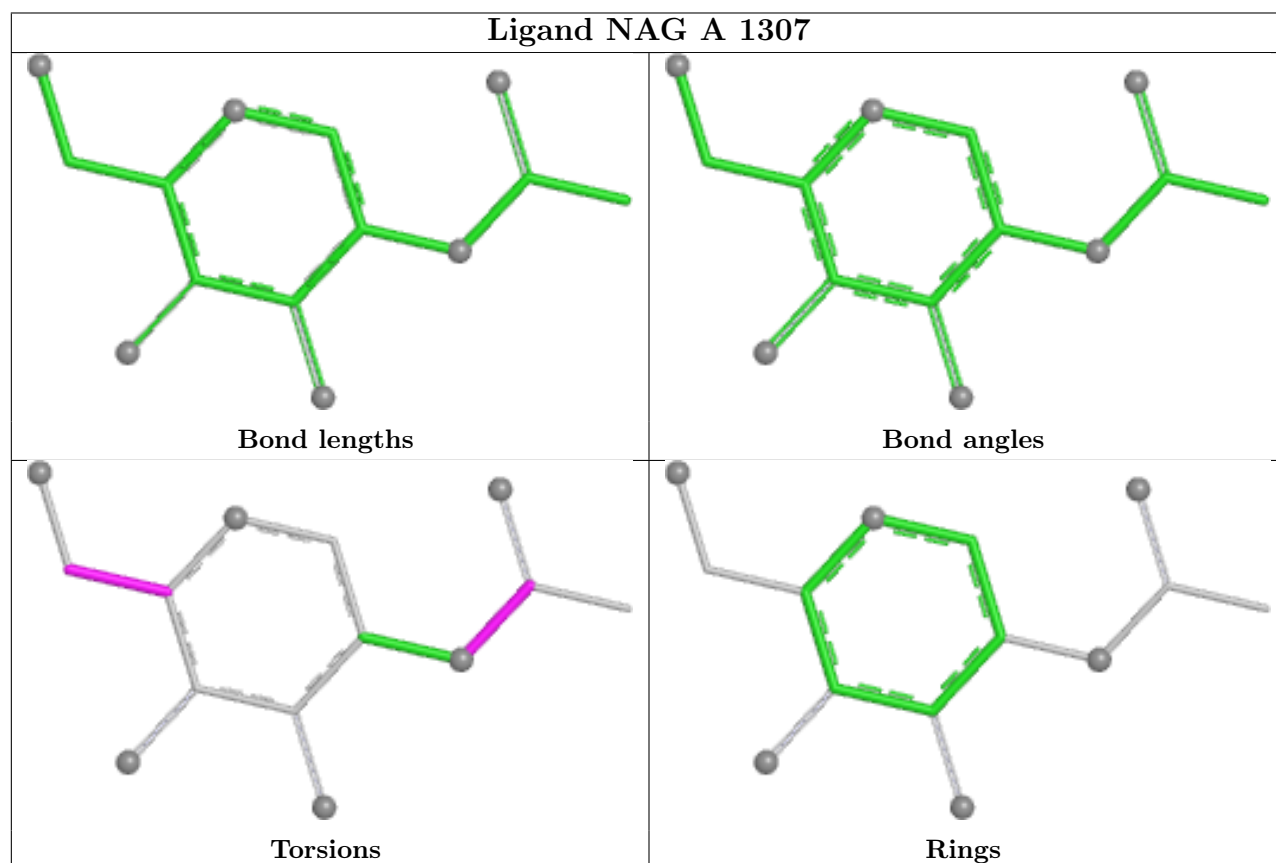
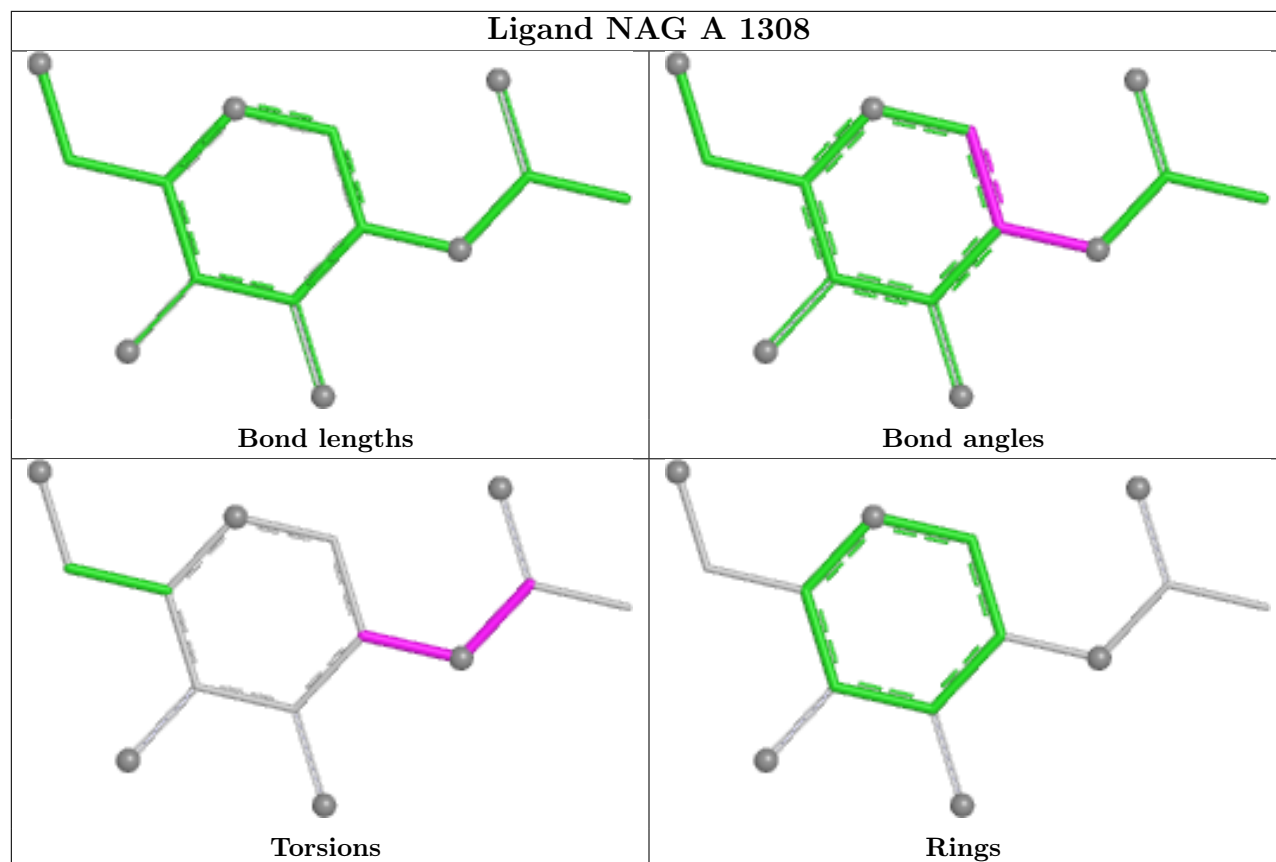


Ligand NAG C 1303

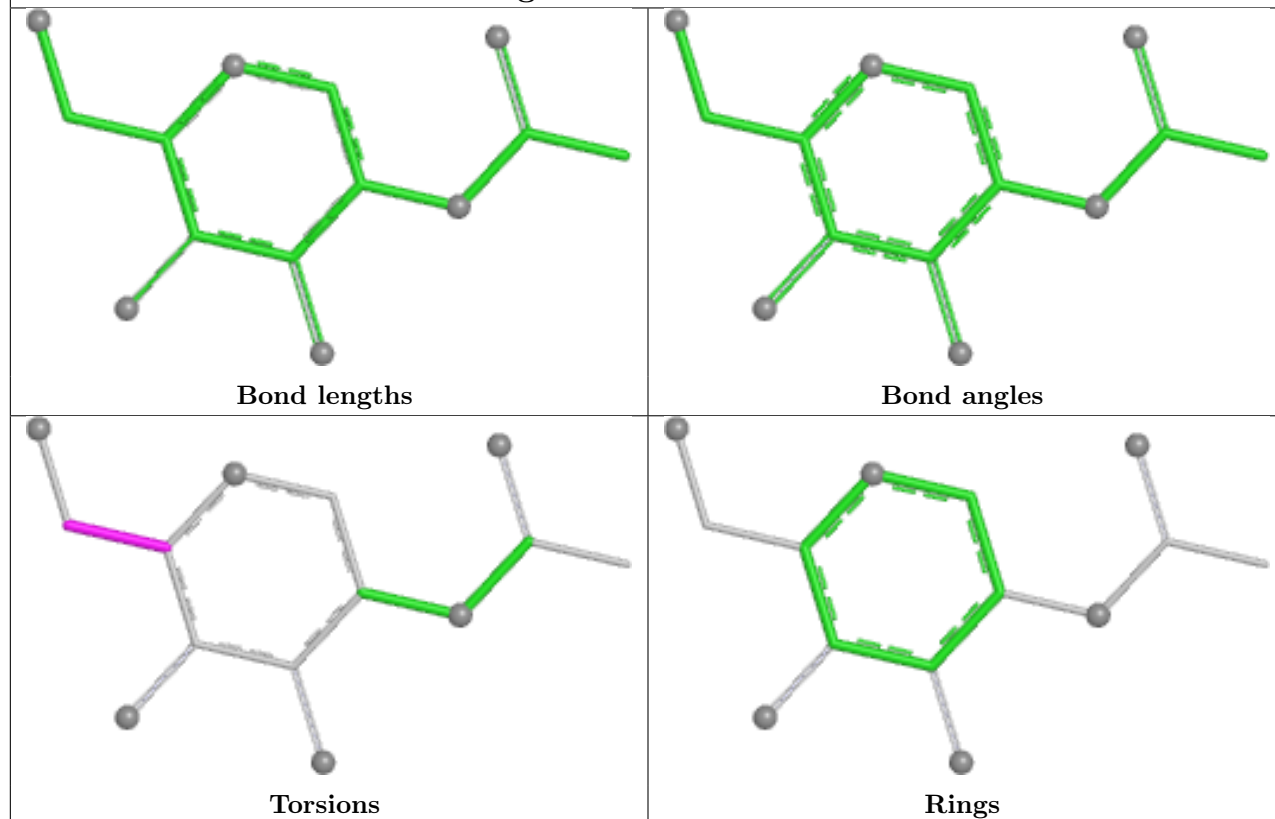


Ligand NAG A 1305

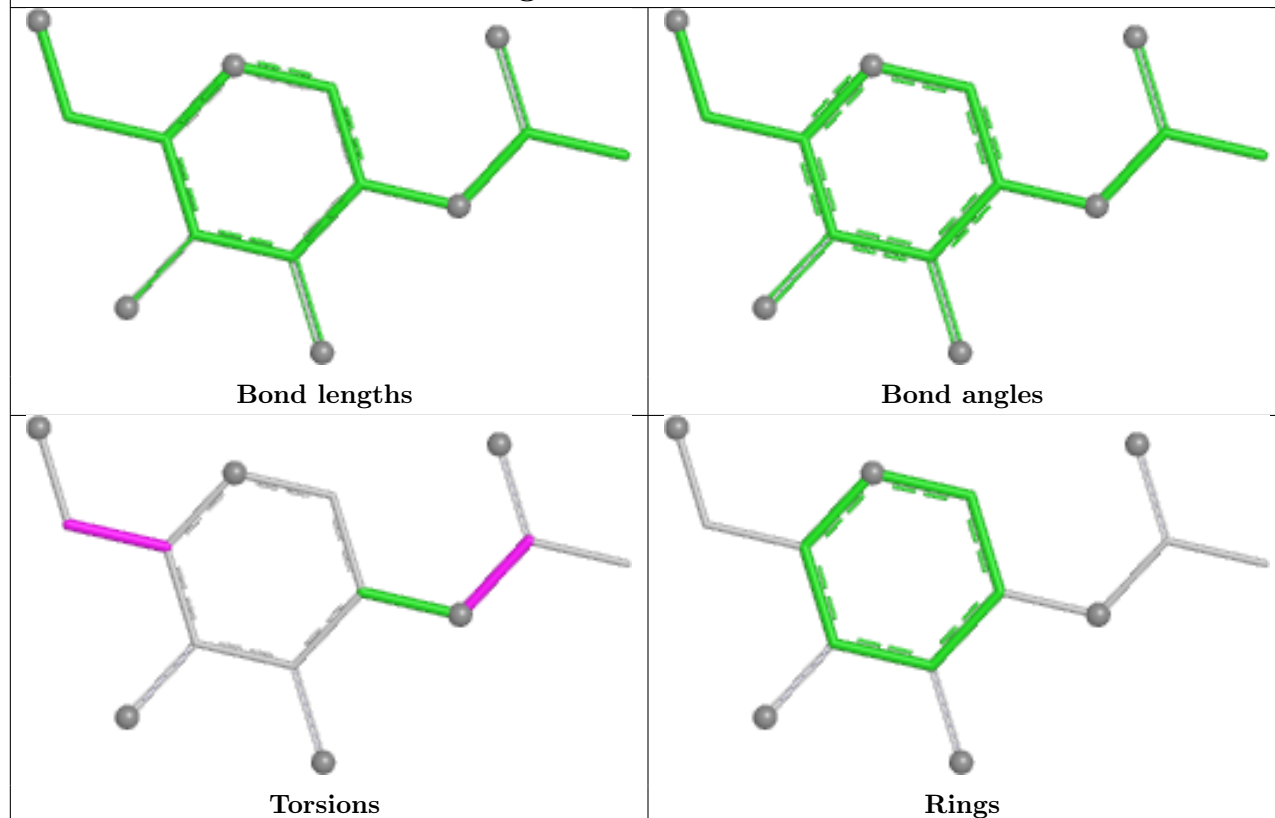




Ligand NAG A 1301



Ligand NAG C 1302



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.

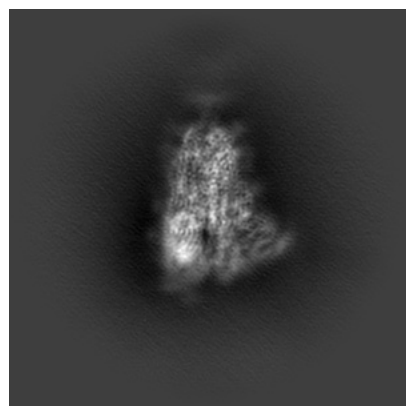
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-38937. These allow visual inspection of the internal detail of the map and identification of artifacts.

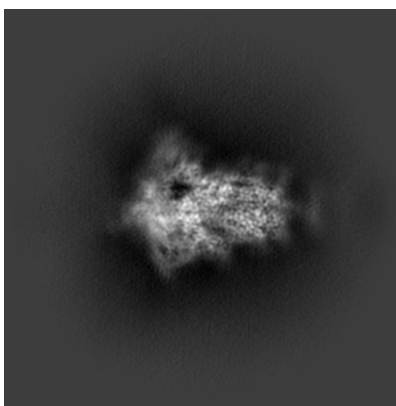
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

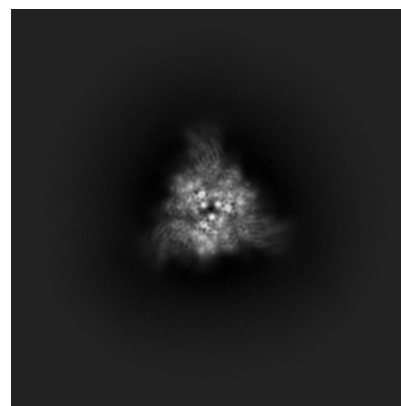
6.1.1 Primary map



X

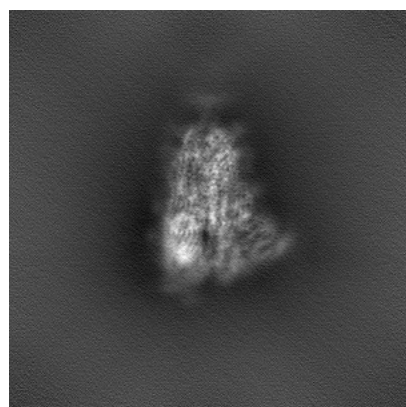


Y

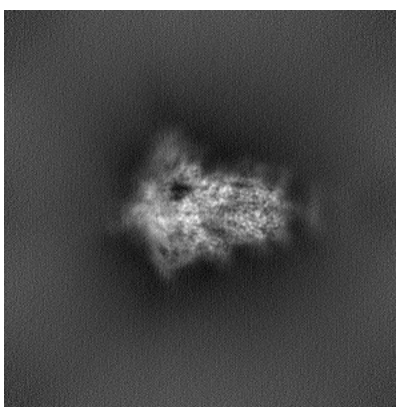


Z

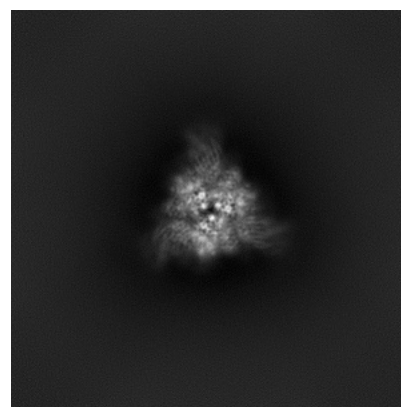
6.1.2 Raw map



X



Y

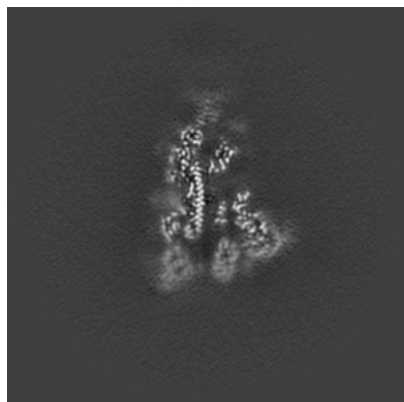


Z

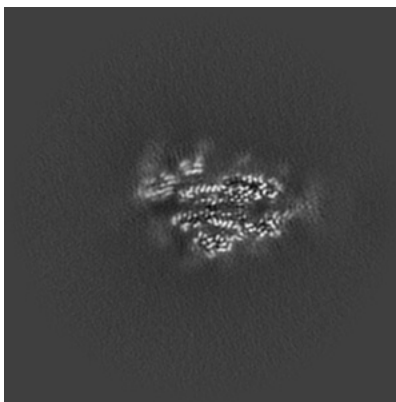
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

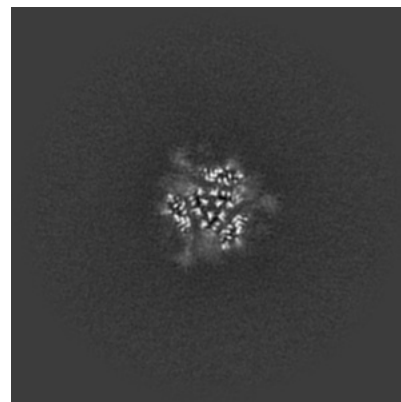
6.2.1 Primary map



X Index: 230

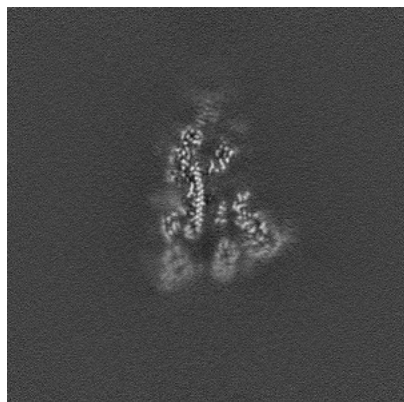


Y Index: 230

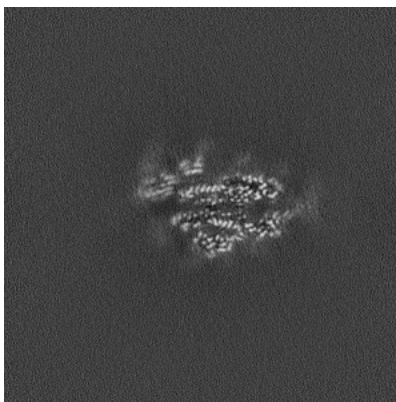


Z Index: 230

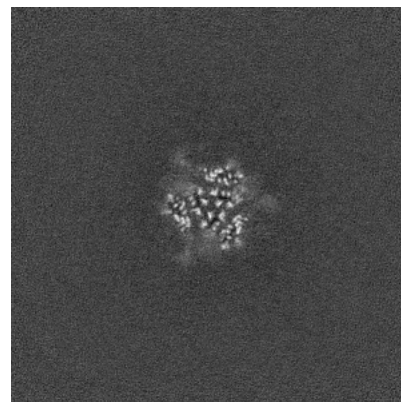
6.2.2 Raw map



X Index: 230



Y Index: 230

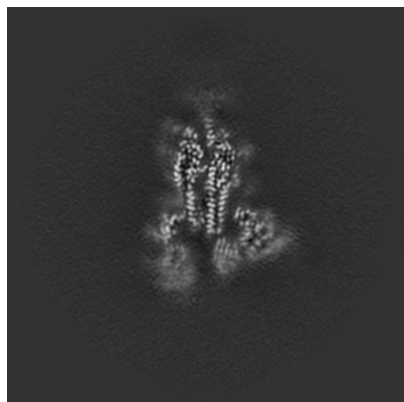


Z Index: 230

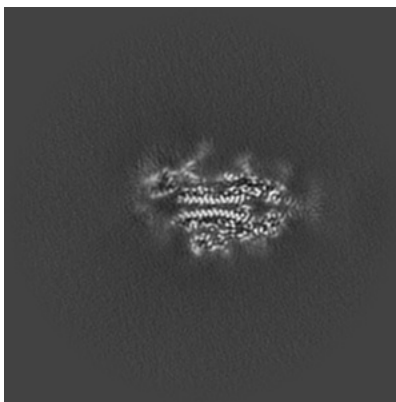
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

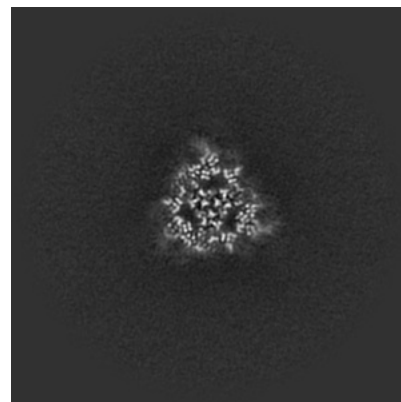
6.3.1 Primary map



X Index: 220

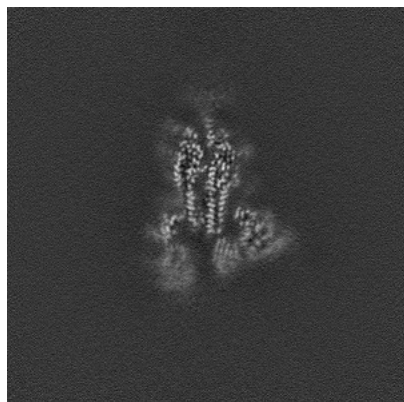


Y Index: 235

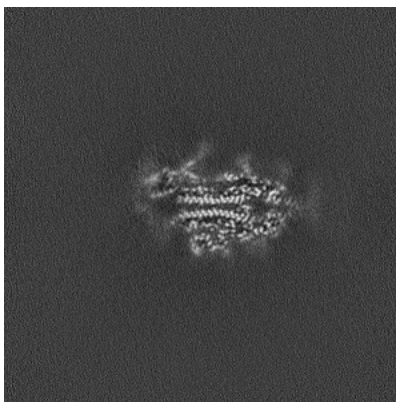


Z Index: 221

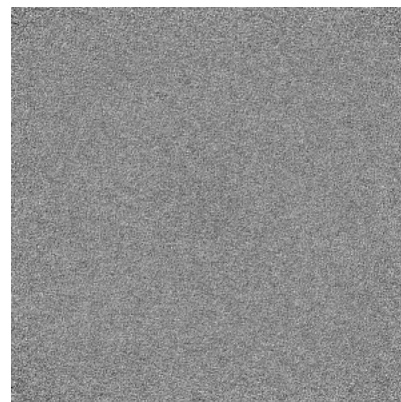
6.3.2 Raw map



X Index: 220



Y Index: 235

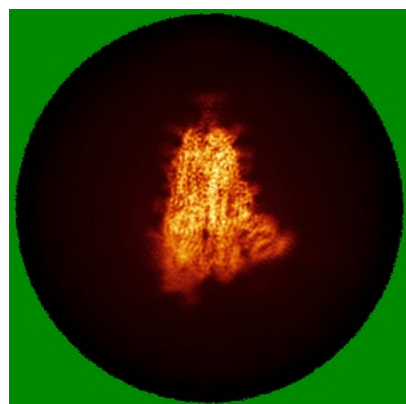


Z Index: 0

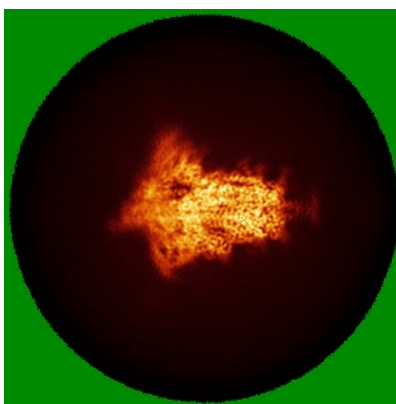
The images above show the largest variance slices of the map in three orthogonal directions.

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

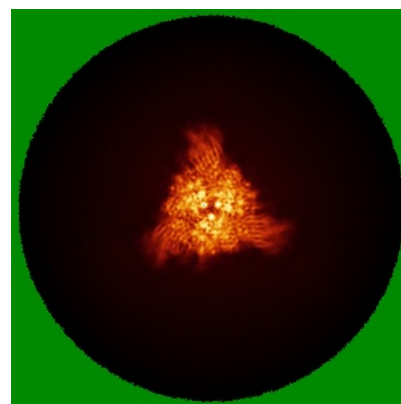
6.4.1 Primary map



X

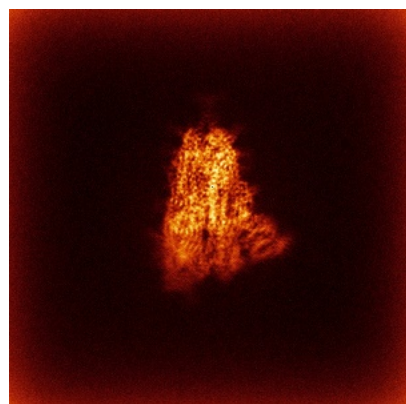


Y

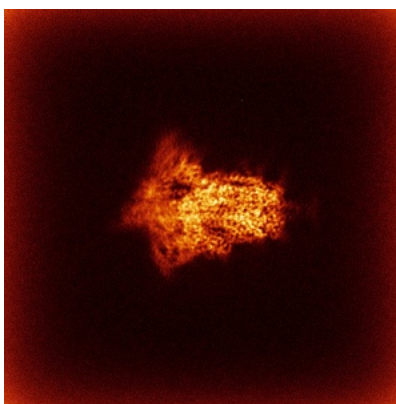


Z

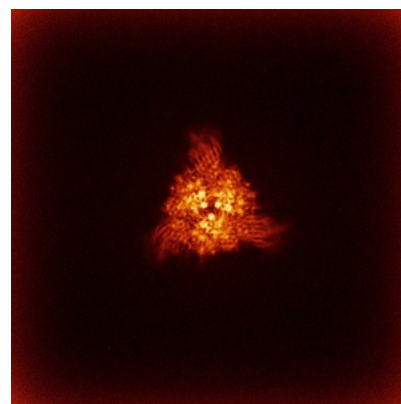
6.4.2 Raw map



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.

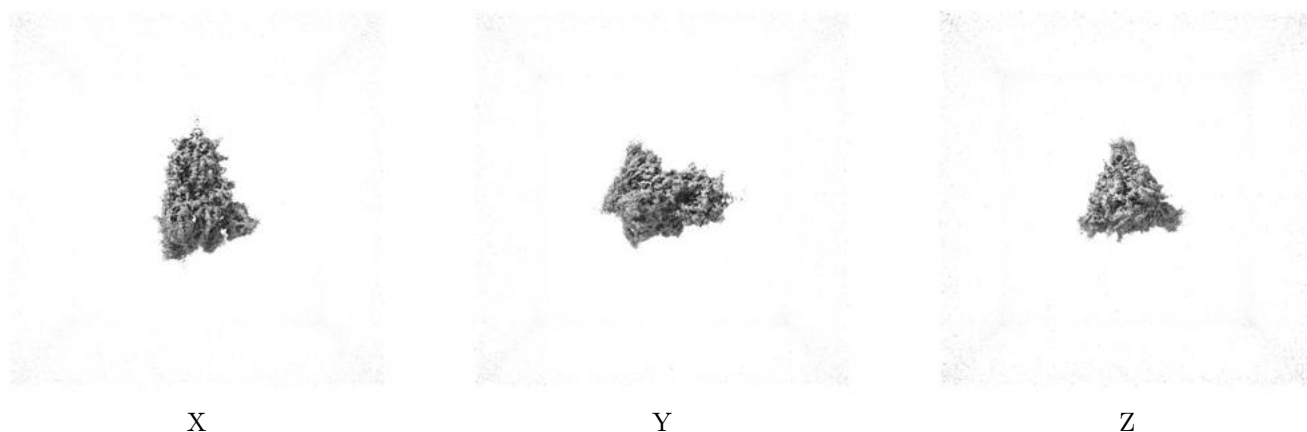
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.2. 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.

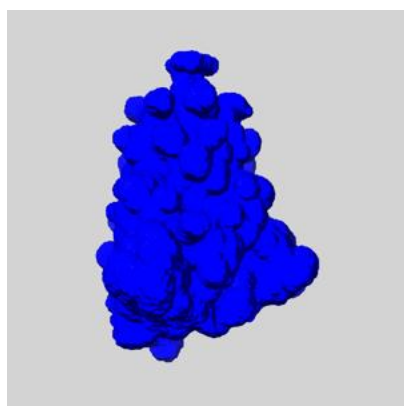
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

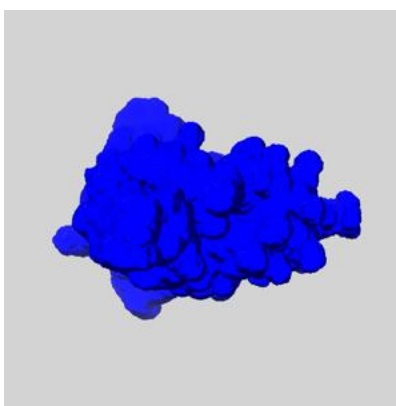
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

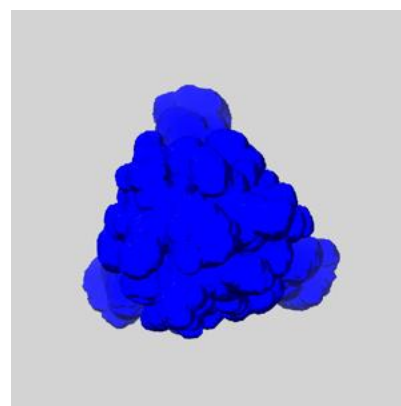
6.6.1 emd_38937_msk_1.map [i](#)



X



Y

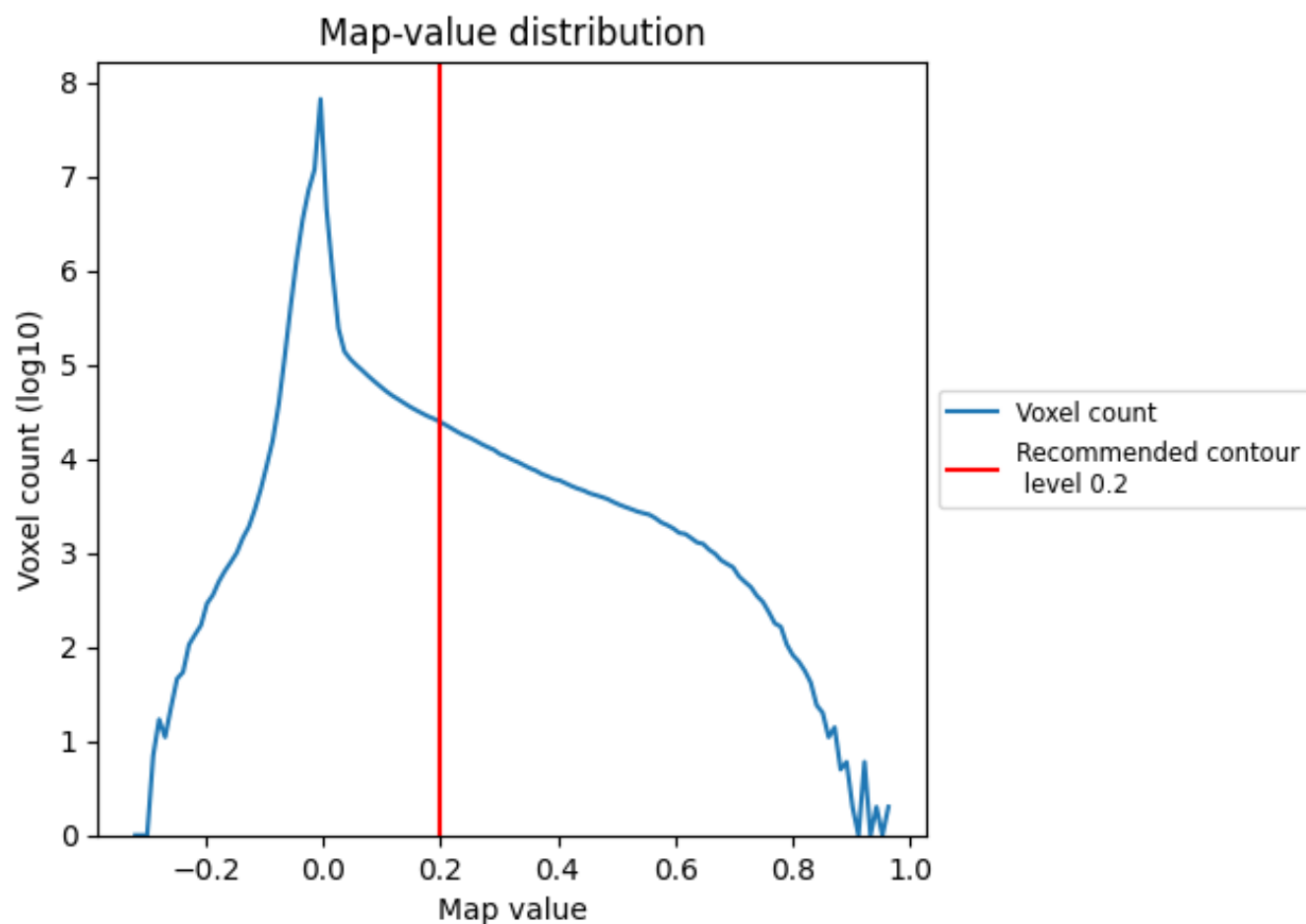


Z

7 Map analysis [i](#)

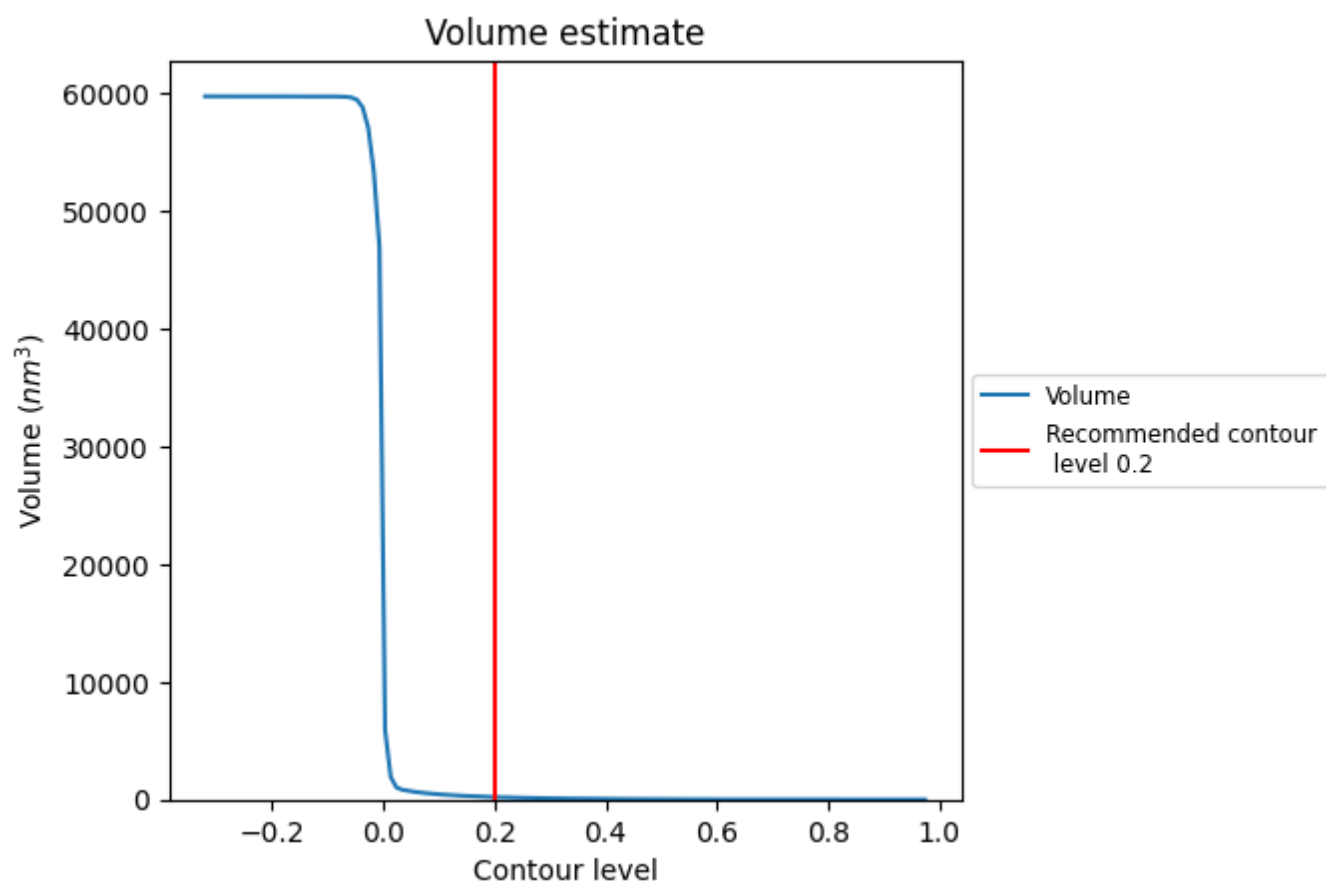
This section contains the results of statistical analysis of the map.

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.

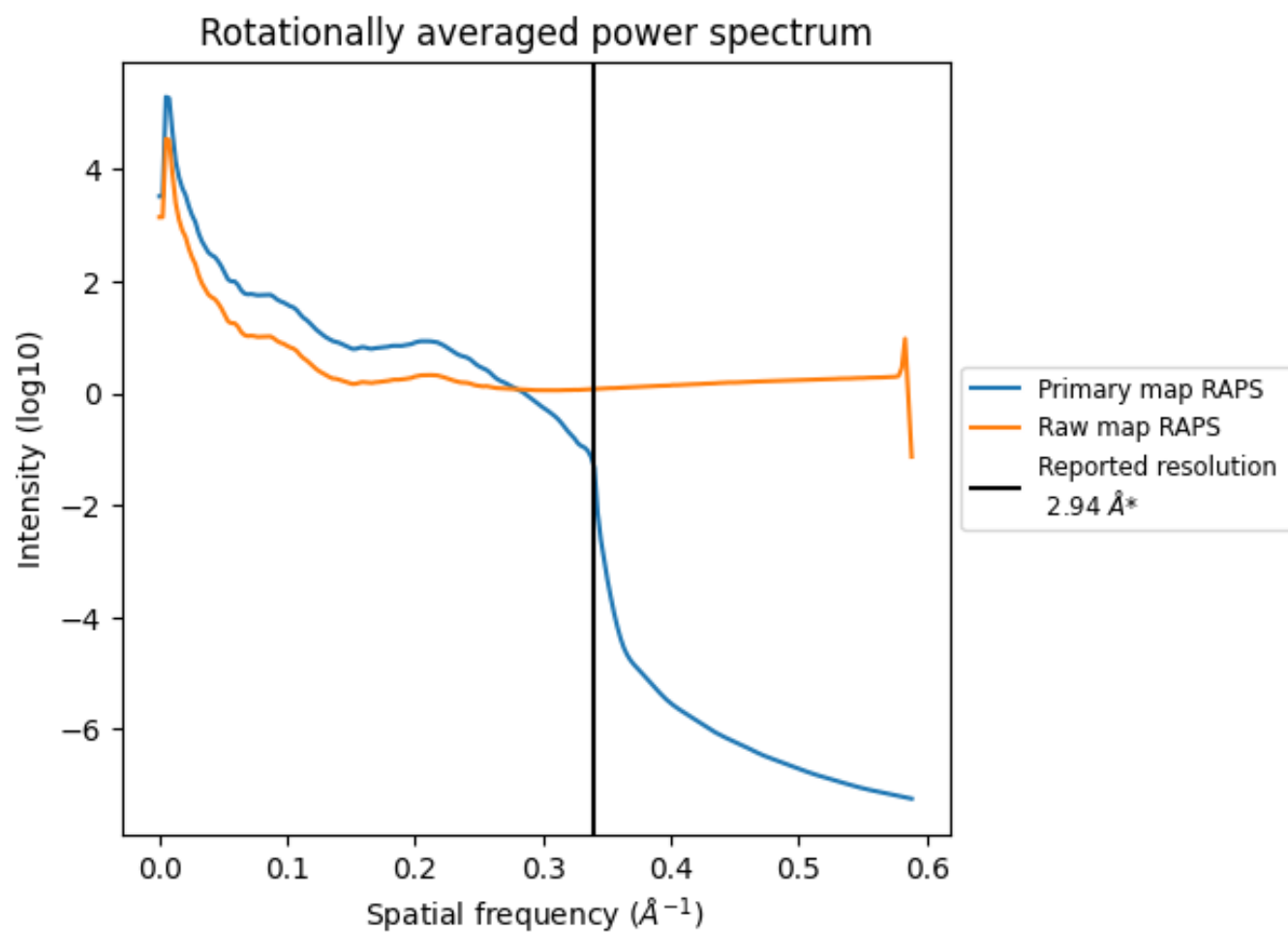
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 212 nm³; this corresponds to an approximate mass of 192 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 ⓘ

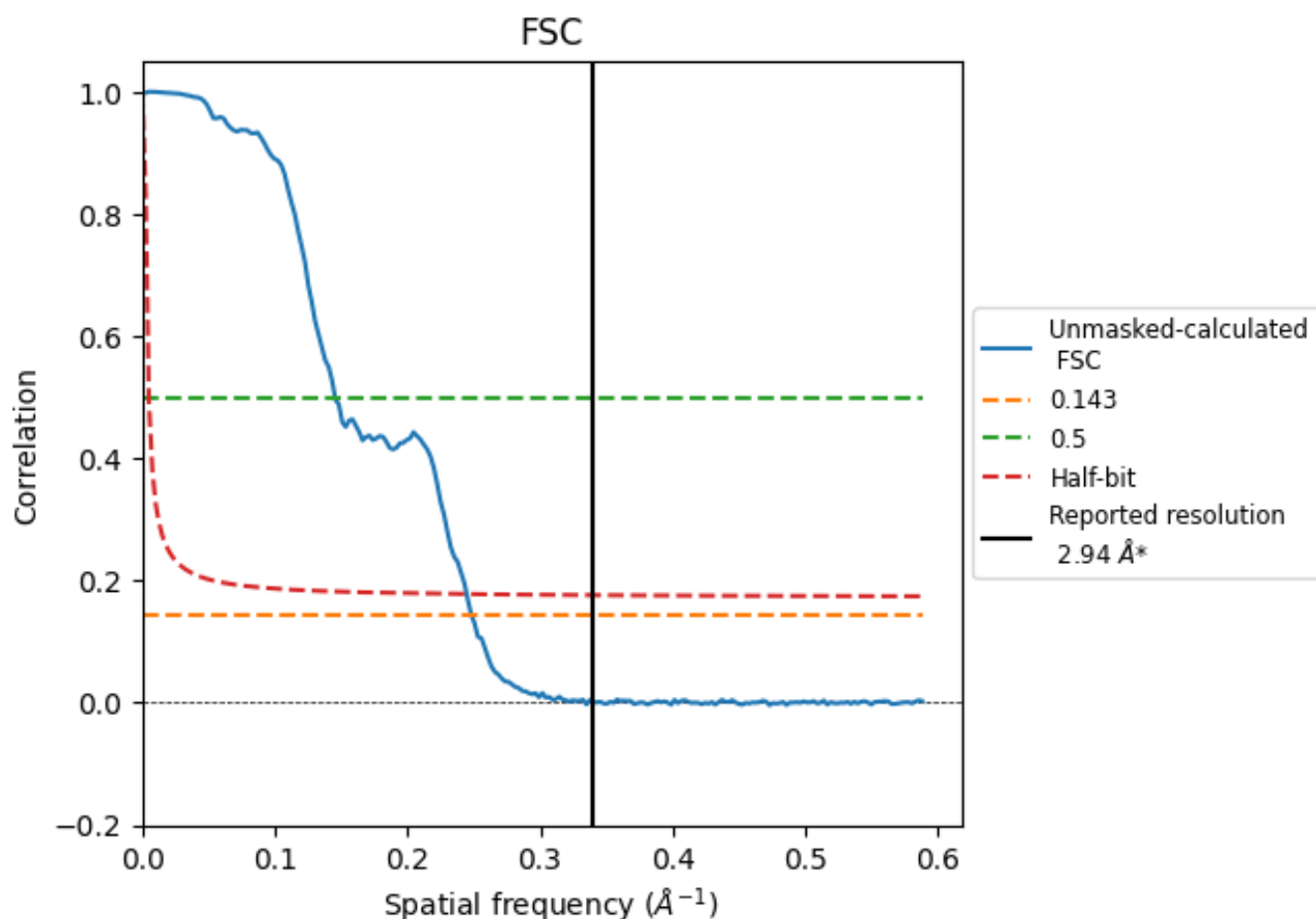


*Reported resolution corresponds to spatial frequency of $0.340 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.340 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.94	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.03	6.88	4.09

*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 4.03 differs from the reported value 2.94 by more than 10 %

9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-38937 and PDB model 8Y5J. Per-residue inclusion information can be found in [section 3](#) on [page 16](#).

9.1 Map-model overlay [i](#)



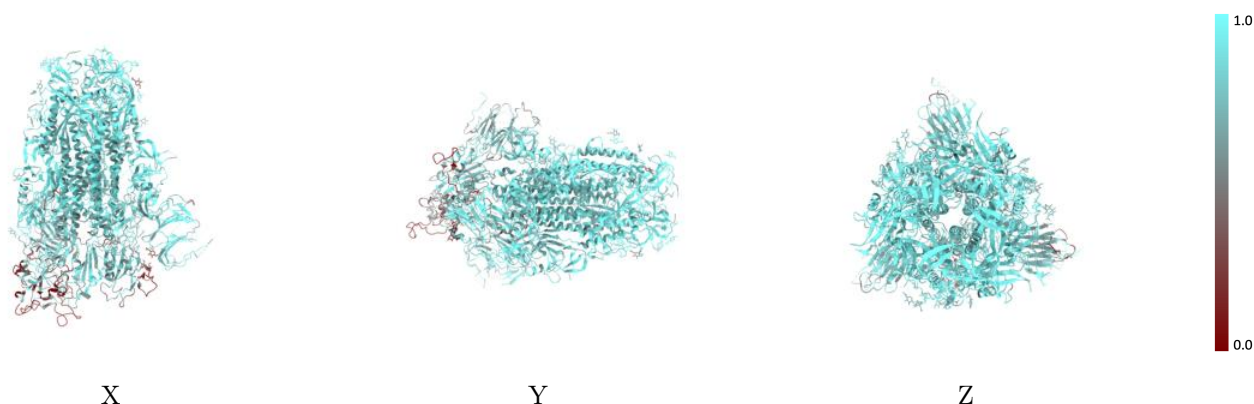
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



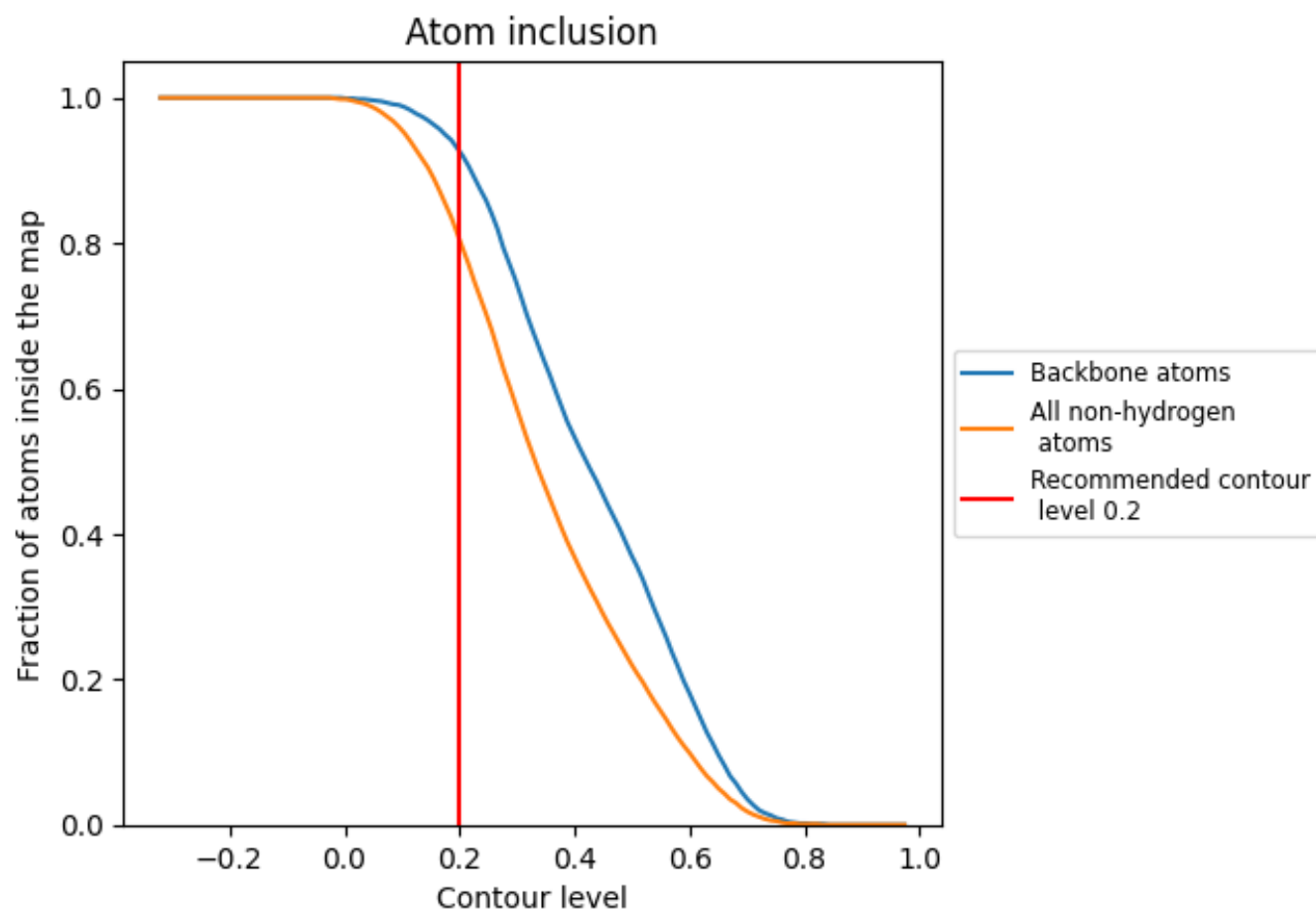
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).

9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 80% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8050	0.4290
A	0.8370	0.4480
B	0.7580	0.3900
C	0.8240	0.4480
D	0.8210	0.4290
E	0.8210	0.4280
F	0.7860	0.4110
G	0.2050	0.2760

1.0

0.0

<0.0