Full wwPDB EM Validation Report ⓘ

Jun 26, 2025 – 09:36 AM JST

PDB ID	:	8H01 / pdb_00008h01
EMDB ID	:	EMD-34408
Title	:	SARS-CoV-2 Omicron BA.1 Spike glycoprotein in complex with rabbit monoclonal antibody 1H1 Fab in class 2 conformation
Authors	:	Guo, H.; Gao, Y.; Lu, Y.; Yang, H.; Ji, X.
Deposited on	:	2022-09-27
Resolution	:	3.70 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

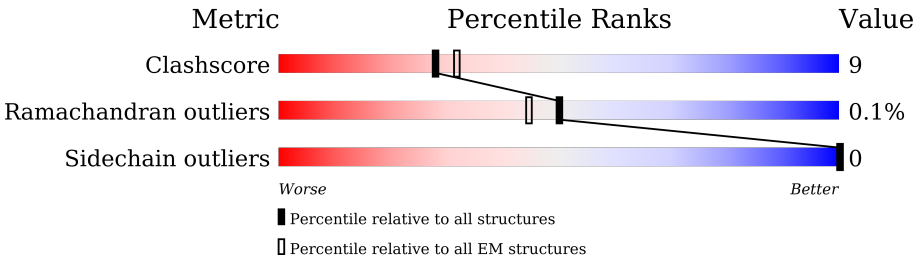
EMDB validation analysis	:	0.0.1.dev118
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.0rc1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.44

1 Overall quality at a glance

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

The reported resolution of this entry is 3.70 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	1286	63% 17% 20%
1	B	1286	63% 17% 20%
1	C	1286	65% 16% 20%
2	D	111	7% 67% 32%
2	F	111	42% 76% 24%
2	H	111	36% 70% 30%
3	E	122	13% 66% 34%
3	G	122	43% 66% 34%

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Mol	Chain	Length	Quality of chain
3	I	122	43% 70% 30%
4	J	2	100%
4	K	2	100%
4	L	2	100%
4	M	2	100%
4	N	2	100%
4	O	2	100%
4	P	2	100%
4	Q	2	50% 50% 50%
4	R	2	100%
4	S	2	100%
4	T	2	100%
4	U	2	100%
4	V	2	50% 50% 50%
4	W	2	100%
4	X	2	100%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 30144 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	1034	Total	C	N	O	S	0	0
			8118	5203	1351	1527	37		
1	B	1035	Total	C	N	O	S	0	0
			8119	5204	1352	1526	37		
1	C	1035	Total	C	N	O	S	0	0
			8113	5201	1348	1527	37		

There are 381 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	VAL	ALA	variant	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	95	ILE	THR	variant	UNP P0DTC2
A	?	-	GLY	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	142	ASP	TYR	variant	UNP P0DTC2
A	210B	ILE	LEU	variant	UNP P0DTC2
A	210E	GLU	-	insertion	UNP P0DTC2
A	210F	PRO	-	insertion	UNP P0DTC2
A	210G	GLU	-	insertion	UNP P0DTC2
A	339	ASP	GLY	variant	UNP P0DTC2
A	371	LEU	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	446	SER	GLY	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	484	ALA	GLU	variant	UNP P0DTC2
A	493	ARG	GLN	variant	UNP P0DTC2
A	496	SER	GLY	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2
A	547	LYS	THR	variant	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	655	TYR	HIS	variant	UNP P0DTC2
A	679	LYS	ASN	variant	UNP P0DTC2
A	681	HIS	PRO	variant	UNP P0DTC2
A	682	GLY	ARG	engineered mutation	UNP P0DTC2
A	683	SER	ARG	engineered mutation	UNP P0DTC2
A	685	SER	ARG	engineered mutation	UNP P0DTC2
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	817	PRO	PHE	engineered mutation	UNP P0DTC2
A	856	LYS	ASN	variant	UNP P0DTC2
A	892	PRO	ALA	engineered mutation	UNP P0DTC2
A	899	PRO	ALA	engineered mutation	UNP P0DTC2
A	942	PRO	ALA	engineered mutation	UNP P0DTC2
A	954	HIS	GLN	variant	UNP P0DTC2
A	969	LYS	ASN	variant	UNP P0DTC2
A	981	PHE	LEU	variant	UNP P0DTC2
A	986	PRO	LYS	engineered mutation	UNP P0DTC2
A	987	PRO	VAL	engineered mutation	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

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1228	GLY	-	expression tag	UNP P0DTC2
A	1229	GLU	-	expression tag	UNP P0DTC2
A	1230	TRP	-	expression tag	UNP P0DTC2
A	1231	VAL	-	expression tag	UNP P0DTC2
A	1232	PHE	-	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	SER	-	expression tag	UNP P0DTC2
A	1239	GLY	-	expression tag	UNP P0DTC2
A	1240	LEU	-	expression tag	UNP P0DTC2
A	1241	GLU	-	expression tag	UNP P0DTC2
A	1242	VAL	-	expression tag	UNP P0DTC2
A	1243	LEU	-	expression tag	UNP P0DTC2
A	1244	PHE	-	expression tag	UNP P0DTC2
A	1245	GLN	-	expression tag	UNP P0DTC2
A	1246	GLY	-	expression tag	UNP P0DTC2
A	1247	PRO	-	expression tag	UNP P0DTC2
A	1248	GLY	-	expression tag	UNP P0DTC2
A	1249	GLY	-	expression tag	UNP P0DTC2
A	1250	TRP	-	expression tag	UNP P0DTC2
A	1251	SER	-	expression tag	UNP P0DTC2
A	1252	HIS	-	expression tag	UNP P0DTC2
A	1253	PRO	-	expression tag	UNP P0DTC2
A	1254	GLN	-	expression tag	UNP P0DTC2
A	1255	PHE	-	expression tag	UNP P0DTC2
A	1256	GLU	-	expression tag	UNP P0DTC2
A	1257	LYS	-	expression tag	UNP P0DTC2
A	1258	GLY	-	expression tag	UNP P0DTC2
A	1259	GLY	-	expression tag	UNP P0DTC2
A	1260	GLY	-	expression tag	UNP P0DTC2
A	1261	SER	-	expression tag	UNP P0DTC2
A	1262	GLY	-	expression tag	UNP P0DTC2
A	1263	GLY	-	expression tag	UNP P0DTC2
A	1264	GLY	-	expression tag	UNP P0DTC2
A	1265	SER	-	expression tag	UNP P0DTC2
A	1266	GLY	-	expression tag	UNP P0DTC2
A	1267	GLY	-	expression tag	UNP P0DTC2
A	1268	SER	-	expression tag	UNP P0DTC2
A	1269	ALA	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1270	TRP	-	expression tag	UNP P0DTC2
A	1271	SER	-	expression tag	UNP P0DTC2
A	1272	HIS	-	expression tag	UNP P0DTC2
A	1273	PRO	-	expression tag	UNP P0DTC2
A	1274	GLN	-	expression tag	UNP P0DTC2
A	1275	PHE	-	expression tag	UNP P0DTC2
A	1276	GLU	-	expression tag	UNP P0DTC2
A	1277	LYS	-	expression tag	UNP P0DTC2
A	1278	GLY	-	expression tag	UNP P0DTC2
A	1279	GLY	-	expression tag	UNP P0DTC2
A	1280	SER	-	expression tag	UNP P0DTC2
A	1281	HIS	-	expression tag	UNP P0DTC2
A	1282	HIS	-	expression tag	UNP P0DTC2
A	1283	HIS	-	expression tag	UNP P0DTC2
A	1284	HIS	-	expression tag	UNP P0DTC2
A	1285	HIS	-	expression tag	UNP P0DTC2
A	1286	HIS	-	expression tag	UNP P0DTC2
A	1287	HIS	-	expression tag	UNP P0DTC2
A	1288	HIS	-	expression tag	UNP P0DTC2
B	67	VAL	ALA	variant	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	95	ILE	THR	variant	UNP P0DTC2
B	?	-	GLY	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	142	ASP	TYR	variant	UNP P0DTC2
B	210B	ILE	LEU	variant	UNP P0DTC2
B	210E	GLU	-	insertion	UNP P0DTC2
B	210F	PRO	-	insertion	UNP P0DTC2
B	210G	GLU	-	insertion	UNP P0DTC2
B	339	ASP	GLY	variant	UNP P0DTC2
B	371	LEU	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	446	SER	GLY	variant	UNP P0DTC2
B	477	ASN	SER	variant	UNP P0DTC2
B	478	LYS	THR	variant	UNP P0DTC2
B	484	ALA	GLU	variant	UNP P0DTC2
B	493	ARG	GLN	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	496	SER	GLY	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	505	HIS	TYR	variant	UNP P0DTC2
B	547	LYS	THR	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	679	LYS	ASN	variant	UNP P0DTC2
B	681	HIS	PRO	variant	UNP P0DTC2
B	682	GLY	ARG	engineered mutation	UNP P0DTC2
B	683	SER	ARG	engineered mutation	UNP P0DTC2
B	685	SER	ARG	engineered mutation	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	817	PRO	PHE	engineered mutation	UNP P0DTC2
B	856	LYS	ASN	variant	UNP P0DTC2
B	892	PRO	ALA	engineered mutation	UNP P0DTC2
B	899	PRO	ALA	engineered mutation	UNP P0DTC2
B	942	PRO	ALA	engineered mutation	UNP P0DTC2
B	954	HIS	GLN	variant	UNP P0DTC2
B	969	LYS	ASN	variant	UNP P0DTC2
B	981	PHE	LEU	variant	UNP P0DTC2
B	986	PRO	LYS	engineered mutation	UNP P0DTC2
B	987	PRO	VAL	engineered mutation	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

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1227	ASP	-	expression tag	UNP P0DTC2
B	1228	GLY	-	expression tag	UNP P0DTC2
B	1229	GLU	-	expression tag	UNP P0DTC2
B	1230	TRP	-	expression tag	UNP P0DTC2
B	1231	VAL	-	expression tag	UNP P0DTC2
B	1232	PHE	-	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	SER	-	expression tag	UNP P0DTC2
B	1239	GLY	-	expression tag	UNP P0DTC2
B	1240	LEU	-	expression tag	UNP P0DTC2
B	1241	GLU	-	expression tag	UNP P0DTC2
B	1242	VAL	-	expression tag	UNP P0DTC2
B	1243	LEU	-	expression tag	UNP P0DTC2
B	1244	PHE	-	expression tag	UNP P0DTC2
B	1245	GLN	-	expression tag	UNP P0DTC2
B	1246	GLY	-	expression tag	UNP P0DTC2
B	1247	PRO	-	expression tag	UNP P0DTC2
B	1248	GLY	-	expression tag	UNP P0DTC2
B	1249	GLY	-	expression tag	UNP P0DTC2
B	1250	TRP	-	expression tag	UNP P0DTC2
B	1251	SER	-	expression tag	UNP P0DTC2
B	1252	HIS	-	expression tag	UNP P0DTC2
B	1253	PRO	-	expression tag	UNP P0DTC2
B	1254	GLN	-	expression tag	UNP P0DTC2
B	1255	PHE	-	expression tag	UNP P0DTC2
B	1256	GLU	-	expression tag	UNP P0DTC2
B	1257	LYS	-	expression tag	UNP P0DTC2
B	1258	GLY	-	expression tag	UNP P0DTC2
B	1259	GLY	-	expression tag	UNP P0DTC2
B	1260	GLY	-	expression tag	UNP P0DTC2
B	1261	SER	-	expression tag	UNP P0DTC2
B	1262	GLY	-	expression tag	UNP P0DTC2
B	1263	GLY	-	expression tag	UNP P0DTC2
B	1264	GLY	-	expression tag	UNP P0DTC2
B	1265	SER	-	expression tag	UNP P0DTC2
B	1266	GLY	-	expression tag	UNP P0DTC2
B	1267	GLY	-	expression tag	UNP P0DTC2
B	1268	SER	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1269	ALA	-	expression tag	UNP P0DTC2
B	1270	TRP	-	expression tag	UNP P0DTC2
B	1271	SER	-	expression tag	UNP P0DTC2
B	1272	HIS	-	expression tag	UNP P0DTC2
B	1273	PRO	-	expression tag	UNP P0DTC2
B	1274	GLN	-	expression tag	UNP P0DTC2
B	1275	PHE	-	expression tag	UNP P0DTC2
B	1276	GLU	-	expression tag	UNP P0DTC2
B	1277	LYS	-	expression tag	UNP P0DTC2
B	1278	GLY	-	expression tag	UNP P0DTC2
B	1279	GLY	-	expression tag	UNP P0DTC2
B	1280	SER	-	expression tag	UNP P0DTC2
B	1281	HIS	-	expression tag	UNP P0DTC2
B	1282	HIS	-	expression tag	UNP P0DTC2
B	1283	HIS	-	expression tag	UNP P0DTC2
B	1284	HIS	-	expression tag	UNP P0DTC2
B	1285	HIS	-	expression tag	UNP P0DTC2
B	1286	HIS	-	expression tag	UNP P0DTC2
B	1287	HIS	-	expression tag	UNP P0DTC2
B	1288	HIS	-	expression tag	UNP P0DTC2
C	67	VAL	ALA	variant	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	95	ILE	THR	variant	UNP P0DTC2
C	?	-	GLY	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	142	ASP	TYR	variant	UNP P0DTC2
C	210B	ILE	LEU	variant	UNP P0DTC2
C	210E	GLU	-	insertion	UNP P0DTC2
C	210F	PRO	-	insertion	UNP P0DTC2
C	210G	GLU	-	insertion	UNP P0DTC2
C	339	ASP	GLY	variant	UNP P0DTC2
C	371	LEU	SER	variant	UNP P0DTC2
C	373	PRO	SER	variant	UNP P0DTC2
C	375	PHE	SER	variant	UNP P0DTC2
C	417	ASN	LYS	variant	UNP P0DTC2
C	440	LYS	ASN	variant	UNP P0DTC2
C	446	SER	GLY	variant	UNP P0DTC2
C	477	ASN	SER	variant	UNP P0DTC2
C	478	LYS	THR	variant	UNP P0DTC2
C	484	ALA	GLU	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	493	ARG	GLN	variant	UNP P0DTC2
C	496	SER	GLY	variant	UNP P0DTC2
C	498	ARG	GLN	variant	UNP P0DTC2
C	501	TYR	ASN	variant	UNP P0DTC2
C	505	HIS	TYR	variant	UNP P0DTC2
C	547	LYS	THR	variant	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	655	TYR	HIS	variant	UNP P0DTC2
C	679	LYS	ASN	variant	UNP P0DTC2
C	681	HIS	PRO	variant	UNP P0DTC2
C	682	GLY	ARG	engineered mutation	UNP P0DTC2
C	683	SER	ARG	engineered mutation	UNP P0DTC2
C	685	SER	ARG	engineered mutation	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2
C	817	PRO	PHE	engineered mutation	UNP P0DTC2
C	856	LYS	ASN	variant	UNP P0DTC2
C	892	PRO	ALA	engineered mutation	UNP P0DTC2
C	899	PRO	ALA	engineered mutation	UNP P0DTC2
C	942	PRO	ALA	engineered mutation	UNP P0DTC2
C	954	HIS	GLN	variant	UNP P0DTC2
C	969	LYS	ASN	variant	UNP P0DTC2
C	981	PHE	LEU	variant	UNP P0DTC2
C	986	PRO	LYS	engineered mutation	UNP P0DTC2
C	987	PRO	VAL	engineered mutation	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

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1226	LYS	-	expression tag	UNP P0DTC2
C	1227	ASP	-	expression tag	UNP P0DTC2
C	1228	GLY	-	expression tag	UNP P0DTC2
C	1229	GLU	-	expression tag	UNP P0DTC2
C	1230	TRP	-	expression tag	UNP P0DTC2
C	1231	VAL	-	expression tag	UNP P0DTC2
C	1232	PHE	-	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	SER	-	expression tag	UNP P0DTC2
C	1239	GLY	-	expression tag	UNP P0DTC2
C	1240	LEU	-	expression tag	UNP P0DTC2
C	1241	GLU	-	expression tag	UNP P0DTC2
C	1242	VAL	-	expression tag	UNP P0DTC2
C	1243	LEU	-	expression tag	UNP P0DTC2
C	1244	PHE	-	expression tag	UNP P0DTC2
C	1245	GLN	-	expression tag	UNP P0DTC2
C	1246	GLY	-	expression tag	UNP P0DTC2
C	1247	PRO	-	expression tag	UNP P0DTC2
C	1248	GLY	-	expression tag	UNP P0DTC2
C	1249	GLY	-	expression tag	UNP P0DTC2
C	1250	TRP	-	expression tag	UNP P0DTC2
C	1251	SER	-	expression tag	UNP P0DTC2
C	1252	HIS	-	expression tag	UNP P0DTC2
C	1253	PRO	-	expression tag	UNP P0DTC2
C	1254	GLN	-	expression tag	UNP P0DTC2
C	1255	PHE	-	expression tag	UNP P0DTC2
C	1256	GLU	-	expression tag	UNP P0DTC2
C	1257	LYS	-	expression tag	UNP P0DTC2
C	1258	GLY	-	expression tag	UNP P0DTC2
C	1259	GLY	-	expression tag	UNP P0DTC2
C	1260	GLY	-	expression tag	UNP P0DTC2
C	1261	SER	-	expression tag	UNP P0DTC2
C	1262	GLY	-	expression tag	UNP P0DTC2
C	1263	GLY	-	expression tag	UNP P0DTC2
C	1264	GLY	-	expression tag	UNP P0DTC2
C	1265	SER	-	expression tag	UNP P0DTC2
C	1266	GLY	-	expression tag	UNP P0DTC2
C	1267	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1268	SER	-	expression tag	UNP P0DTC2
C	1269	ALA	-	expression tag	UNP P0DTC2
C	1270	TRP	-	expression tag	UNP P0DTC2
C	1271	SER	-	expression tag	UNP P0DTC2
C	1272	HIS	-	expression tag	UNP P0DTC2
C	1273	PRO	-	expression tag	UNP P0DTC2
C	1274	GLN	-	expression tag	UNP P0DTC2
C	1275	PHE	-	expression tag	UNP P0DTC2
C	1276	GLU	-	expression tag	UNP P0DTC2
C	1277	LYS	-	expression tag	UNP P0DTC2
C	1278	GLY	-	expression tag	UNP P0DTC2
C	1279	GLY	-	expression tag	UNP P0DTC2
C	1280	SER	-	expression tag	UNP P0DTC2
C	1281	HIS	-	expression tag	UNP P0DTC2
C	1282	HIS	-	expression tag	UNP P0DTC2
C	1283	HIS	-	expression tag	UNP P0DTC2
C	1284	HIS	-	expression tag	UNP P0DTC2
C	1285	HIS	-	expression tag	UNP P0DTC2
C	1286	HIS	-	expression tag	UNP P0DTC2
C	1287	HIS	-	expression tag	UNP P0DTC2
C	1288	HIS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called rabbit monoclonal antibody 1H1 Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	111	Total	C	N	O	S	0	0
			824	520	135	165	4		
2	F	111	Total	C	N	O	S	0	0
			824	520	135	165	4		
2	H	111	Total	C	N	O	S	0	0
			824	520	135	165	4		

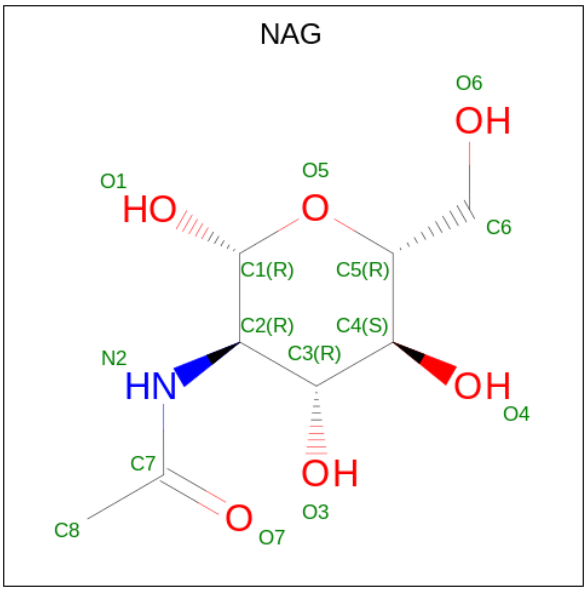
- Molecule 3 is a protein called rabbit monoclonal antibody 1H1 Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	122	Total	C	N	O	S	0	0
			902	568	144	184	6		
3	G	122	Total	C	N	O	S	0	0
			902	568	144	184	6		
3	I	122	Total	C	N	O	S	0	0
			902	568	144	184	6		

- Molecule 4 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
4	O	2	Total	C	N	O	0	0
			28	16	2	10		
4	J	2	Total	C	N	O	0	0
			28	16	2	10		
4	K	2	Total	C	N	O	0	0
			28	16	2	10		
4	L	2	Total	C	N	O	0	0
			28	16	2	10		
4	M	2	Total	C	N	O	0	0
			28	16	2	10		
4	N	2	Total	C	N	O	0	0
			28	16	2	10		
4	P	2	Total	C	N	O	0	0
			28	16	2	10		
4	Q	2	Total	C	N	O	0	0
			28	16	2	10		
4	R	2	Total	C	N	O	0	0
			28	16	2	10		
4	S	2	Total	C	N	O	0	0
			28	16	2	10		
4	T	2	Total	C	N	O	0	0
			28	16	2	10		
4	U	2	Total	C	N	O	0	0
			28	16	2	10		
4	V	2	Total	C	N	O	0	0
			28	16	2	10		
4	W	2	Total	C	N	O	0	0
			28	16	2	10		
4	X	2	Total	C	N	O	0	0
			28	16	2	10		

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



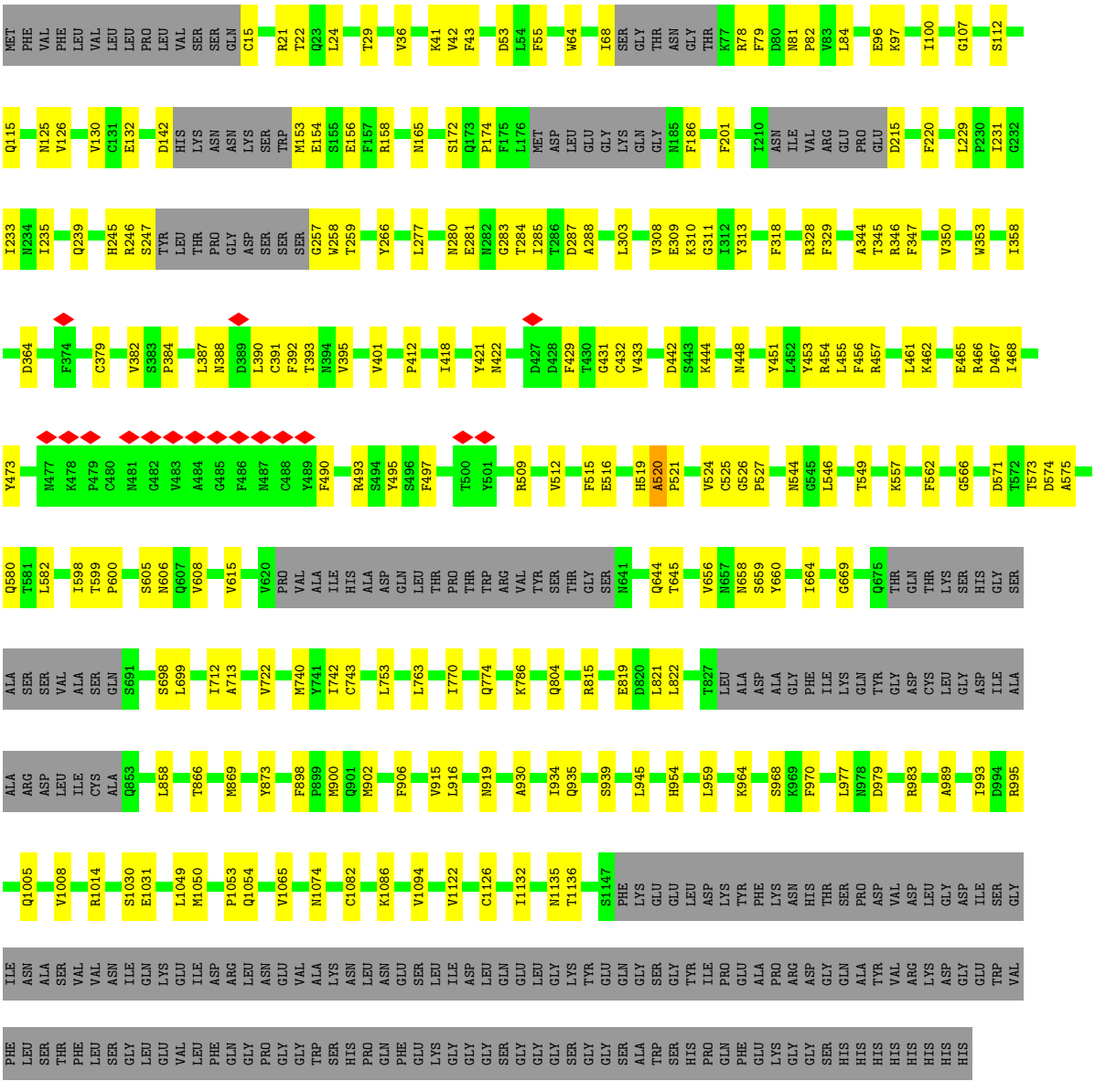
Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	A	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5	B	1	Total	C	N	O	0
			14	8	1	5	
5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

3 Residue-property plots (i)

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.

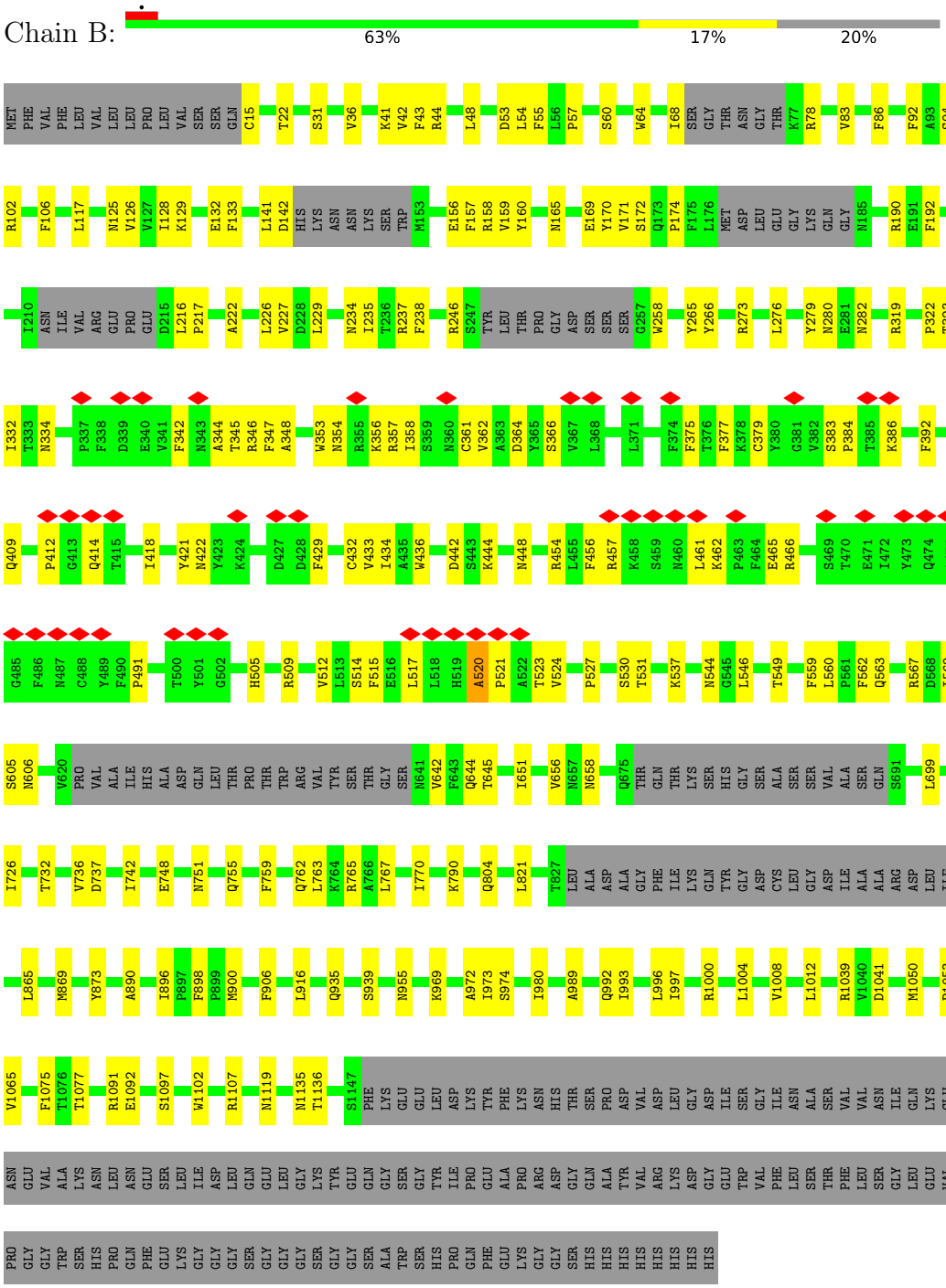
- Molecule 1: Spike glycoprotein

Chain A:



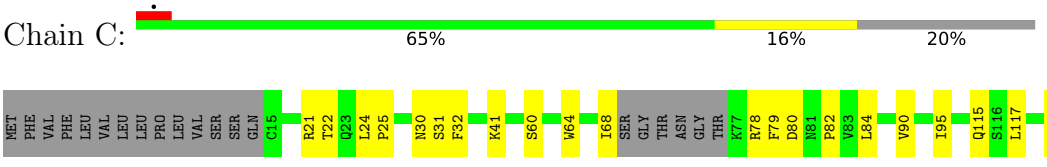
• Molecule 1: Spike glycoprotein

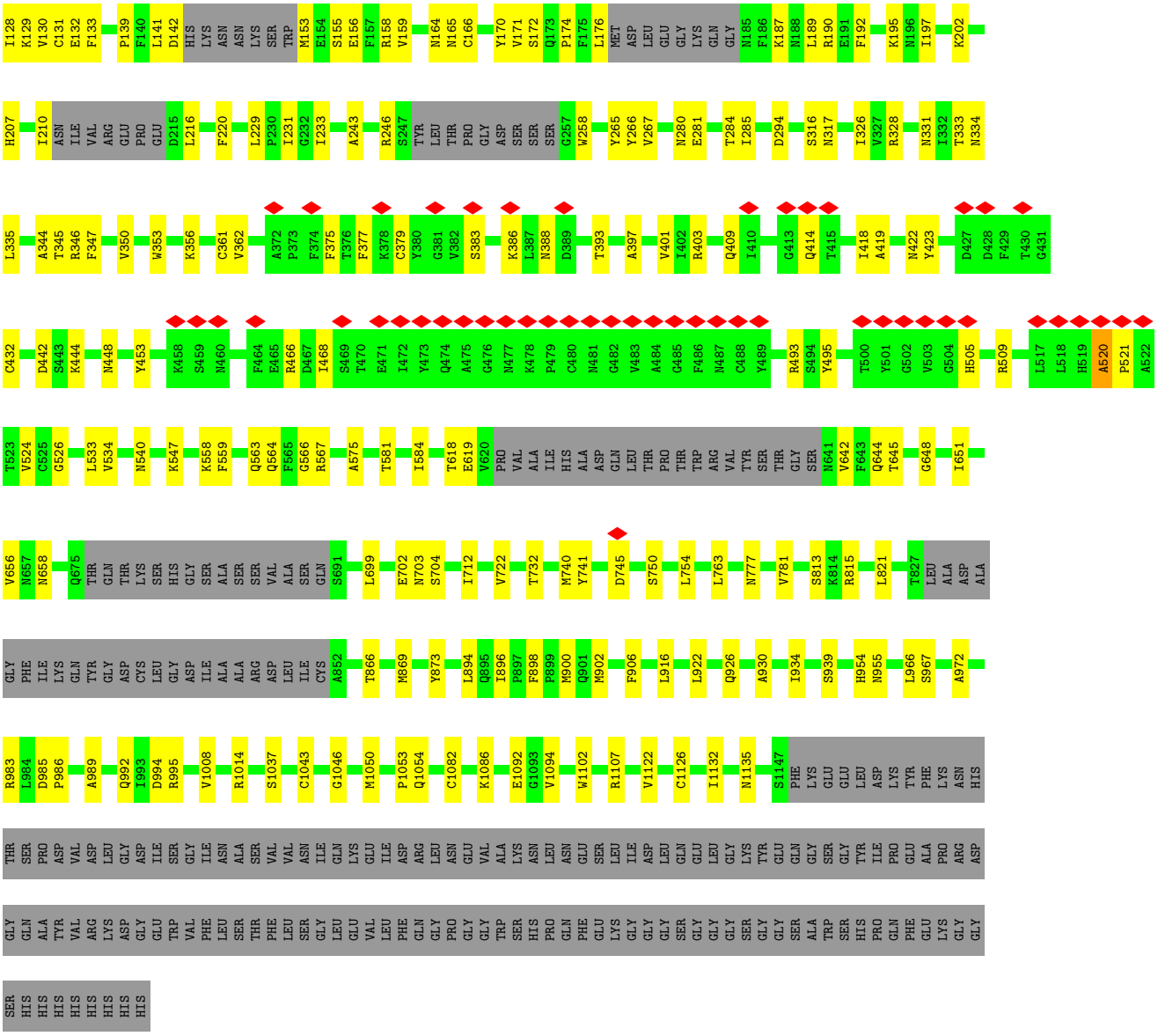
Chain B:



• Molecule 1: Spike glycoprotein

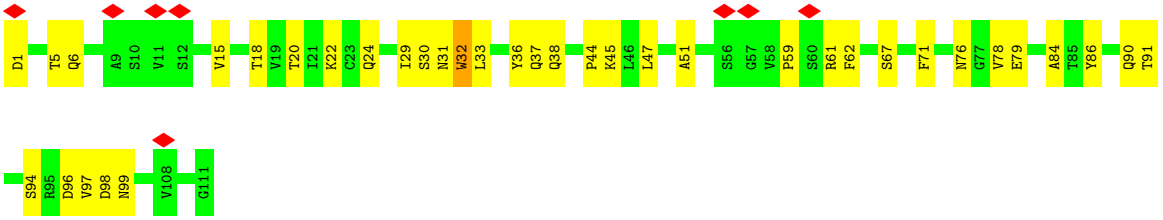
Chain C:





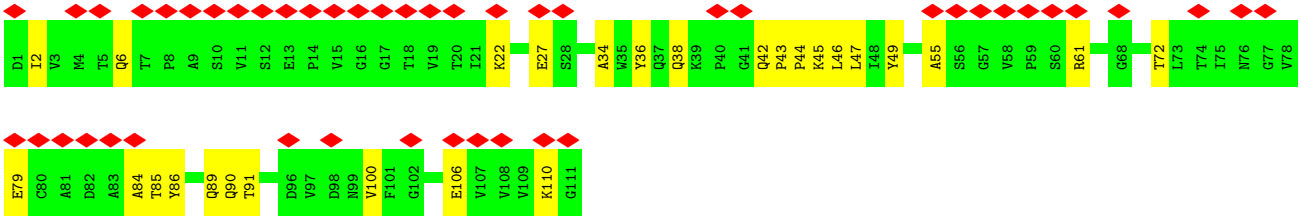
• Molecule 2: rabbit monoclonal antibody 1H1 Fab light chain

Chain D:

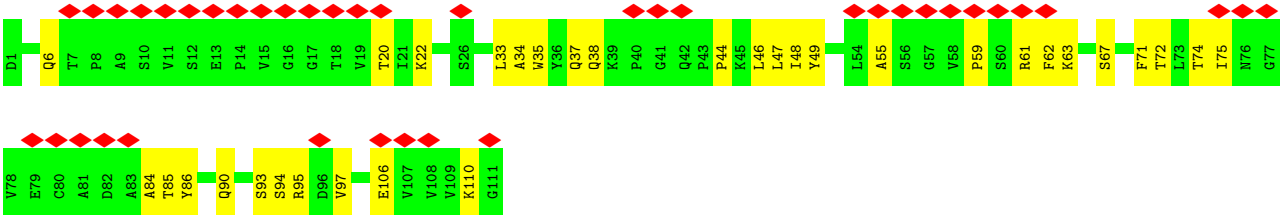


• Molecule 2: rabbit monoclonal antibody 1H1 Fab light chain

Chain F:

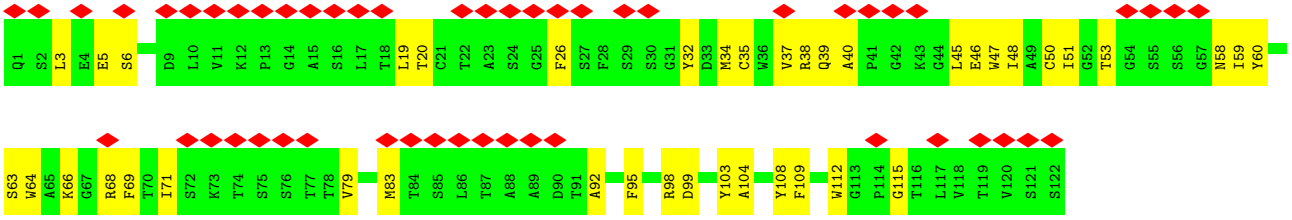


• Molecule 2: rabbit monoclonal antibody 1H1 Fab light chain

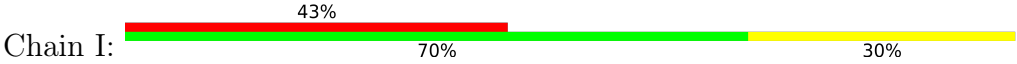


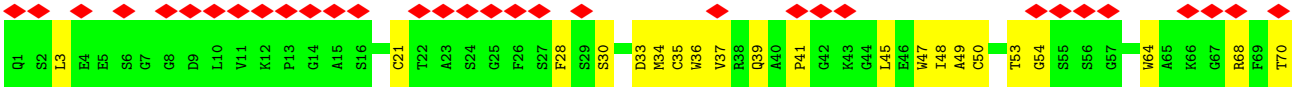
• Molecule 3: rabbit monoclonal antibody 1H1 Fab heavy chain

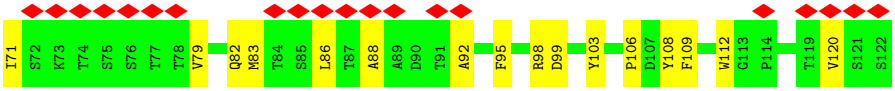
• Molecule 3: rabbit monoclonal antibody 1H1 Fab heavy chain



• Molecule 3: rabbit monoclonal antibody 1H1 Fab heavy chain

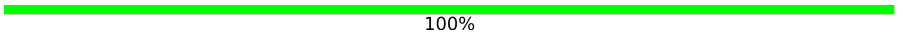






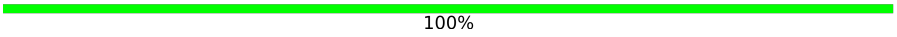
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:



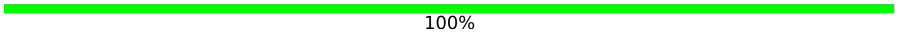
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:



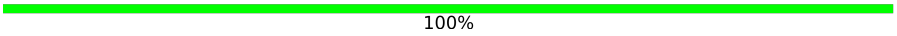
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:



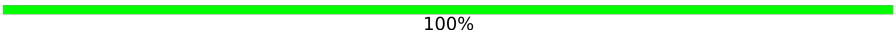
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:



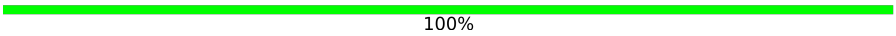
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:



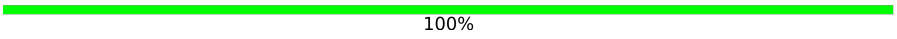
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:



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

Chain P:

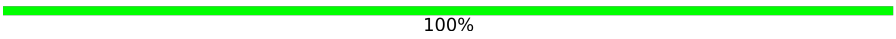


• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:

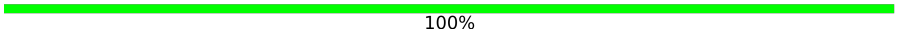
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain R:



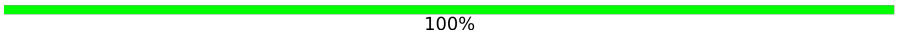
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S:



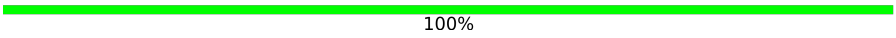
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain T:



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

Chain U:

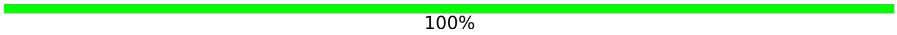


• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain V:

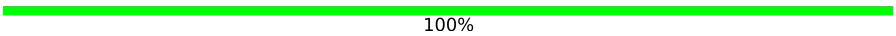
• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain W:



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

Chain X:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	89494	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 ($e^-/Å^2$)	60	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.499	Depositor
Minimum map value	-0.542	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.037	Depositor
Recommended contour level	0.2206	Depositor
Map size (Å)	425.984, 425.984, 425.984	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.832, 0.832, 0.832	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.16	0/8310	0.30	0/11304
1	B	0.15	0/8311	0.30	0/11306
1	C	0.16	0/8305	0.31	0/11299
2	D	0.24	0/842	0.42	0/1148
2	F	0.24	0/842	0.41	0/1148
2	H	0.26	0/842	0.40	0/1148
3	E	0.13	0/924	0.35	0/1259
3	G	0.22	0/924	0.36	0/1259
3	I	0.14	0/924	0.32	0/1259
All	All	0.17	0/30224	0.32	0/41130

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

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	8118	0	7945	150	0
1	B	8119	0	7947	150	0
1	C	8113	0	7932	143	0
2	D	824	0	802	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	824	0	802	23	0
2	H	824	0	802	26	0
3	E	902	0	851	31	0
3	G	902	0	851	35	0
3	I	902	0	851	30	0
4	J	28	0	25	0	0
4	K	28	0	25	0	0
4	L	28	0	25	0	0
4	M	28	0	25	0	0
4	N	28	0	25	0	0
4	O	28	0	25	0	0
4	P	28	0	25	0	0
4	Q	28	0	25	0	0
4	R	28	0	25	0	0
4	S	28	0	25	0	0
4	T	28	0	25	0	0
4	U	28	0	25	0	0
4	V	28	0	25	1	0
4	W	28	0	25	0	0
4	X	28	0	25	0	0
5	A	70	0	65	1	0
5	B	56	0	52	2	0
5	C	70	0	65	0	0
All	All	30144	0	29340	557	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 9.

All (557) 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)
2:F:36:TYR:HB3	2:F:44:PRO:HB2	1.47	0.95
1:A:142:ASP:HB2	1:A:156:GLU:HB3	1.60	0.83
1:A:42:VAL:HG21	1:B:567:ARG:HH21	1.53	0.73
1:B:356:LYS:HB3	1:B:397:ALA:HB3	1.71	0.72
1:B:142:ASP:HB2	1:B:156:GLU:HB3	1.72	0.72
1:C:142:ASP:HB2	1:C:156:GLU:HB3	1.70	0.72
2:H:90:GLN:HE22	2:H:97:VAL:HG21	1.55	0.71
1:A:743:CYS:HB2	1:A:977:LEU:HD22	1.74	0.69
2:H:44:PRO:HG3	3:I:45:LEU:HD11	1.75	0.69
3:I:34:MET:HE2	3:I:79:VAL:HG11	1.73	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:22:LYS:HG3	2:F:72:THR:HG22	1.76	0.68
1:B:353:TRP:O	1:B:466:ARG:NH2	2.27	0.68
1:A:353:TRP:O	1:A:466:ARG:NH1	2.28	0.67
2:D:33:LEU:HB3	2:D:51:ALA:HB2	1.75	0.67
1:B:126:VAL:HG13	1:B:174:PRO:HA	1.77	0.67
1:B:763:LEU:HD13	1:B:1004:LEU:HD22	1.77	0.67
1:C:1053:PRO:O	1:C:1054:GLN:NE2	2.27	0.67
3:G:6:SER:HB3	3:G:20:THR:HB	1.77	0.67
1:C:22:THR:O	1:C:78:ARG:NH1	2.28	0.67
1:C:126:VAL:HG13	1:C:174:PRO:HA	1.78	0.66
2:F:45:LYS:HE2	2:F:47:LEU:HD21	1.77	0.66
1:C:195:LYS:HB2	1:C:202:LYS:HB2	1.78	0.66
1:C:229:LEU:HB3	1:C:231:ILE:HG12	1.77	0.66
1:A:979:ASP:HB3	1:A:983:ARG:HE	1.59	0.66
1:C:345:THR:OG1	2:H:95:ARG:NH2	2.29	0.66
1:A:519:HIS:ND1	1:A:544:ASN:O	2.29	0.66
1:B:656:VAL:HG12	1:B:658:ASN:H	1.60	0.66
1:B:737:ASP:OD2	1:C:317:ASN:ND2	2.29	0.65
1:A:303:LEU:HD22	1:A:308:VAL:HG12	1.79	0.65
1:A:442:ASP:O	1:A:448:ASN:ND2	2.29	0.65
1:C:644:GLN:NE2	1:C:645:THR:O	2.30	0.65
1:C:902:MET:HE1	1:C:1050:MET:HE2	1.79	0.64
1:A:154:GLU:OE2	1:A:245:HIS:ND1	2.30	0.64
1:B:43:PHE:HA	1:C:563:GLN:HE21	1.63	0.64
1:C:153:MET:SD	1:C:155:SER:OG	2.56	0.64
1:B:129:LYS:HG2	1:B:169:GLU:HG3	1.79	0.64
3:E:91:THR:HB	3:E:120:VAL:HG12	1.80	0.64
1:A:869:MET:HG2	1:B:699:LEU:HD11	1.79	0.64
1:A:656:VAL:HG12	1:A:658:ASN:H	1.63	0.63
3:E:2:SER:HB3	3:E:24:SER:HB3	1.79	0.63
1:A:126:VAL:HG23	1:A:174:PRO:HA	1.79	0.63
1:B:396:TYR:HB2	1:B:514:SER:HB3	1.79	0.63
1:C:156:GLU:OE2	1:C:158:ARG:NH1	2.32	0.63
1:A:392:PHE:N	1:A:524:VAL:O	2.30	0.63
1:B:345:THR:O	1:B:509:ARG:NH2	2.30	0.63
1:B:762:GLN:OE1	1:B:765:ARG:NH2	2.31	0.63
1:A:468:ILE:HB	2:D:67:SER:HB3	1.79	0.63
2:F:46:LEU:HD21	2:F:49:TYR:HB3	1.79	0.63
1:A:804:GLN:NE2	1:A:935:GLN:OE1	2.27	0.62
1:B:379:CYS:HB2	1:B:432:CYS:HA	1.79	0.62
1:B:1053:PRO:O	1:B:1054:GLN:NE2	2.29	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:PHE:HB3	1:B:192:PHE:HB2	1.81	0.62
1:B:1092:GLU:O	1:B:1107:ARG:NE	2.31	0.62
1:C:31:SER:HB3	1:C:60:SER:H	1.63	0.62
1:A:379:CYS:HB2	1:A:432:CYS:HA	1.80	0.62
1:A:21:ARG:HG2	1:A:79:PHE:HB2	1.82	0.62
1:C:21:ARG:HB3	1:C:79:PHE:HB2	1.82	0.62
1:B:442:ASP:O	1:B:448:ASN:ND2	2.33	0.62
1:C:821:LEU:HD11	1:C:939:SER:HB2	1.82	0.62
1:B:201:PHE:HB3	1:B:229:LEU:HB2	1.80	0.61
1:A:328:ARG:HH21	1:A:580:GLN:HB2	1.65	0.61
3:E:4:GLU:HA	3:E:114:PRO:HD2	1.82	0.61
1:A:453:TYR:HB3	1:A:495:TYR:HE1	1.66	0.61
1:A:742:ILE:HG21	1:A:753:LEU:HD22	1.82	0.61
1:A:1053:PRO:O	1:A:1054:GLN:NE2	2.30	0.61
1:A:346:ARG:HE	3:E:103:TYR:HB3	1.65	0.61
1:C:197:ILE:H	1:C:202:LYS:HZ1	1.49	0.61
1:A:669:GLY:HA3	1:C:869:MET:HE1	1.83	0.61
2:F:34:ALA:HB2	3:G:108:TYR:HD1	1.65	0.61
1:A:644:GLN:NE2	1:A:645:THR:O	2.34	0.60
1:B:328:ARG:NH1	1:B:531:THR:O	2.34	0.60
1:C:972:ALA:HB2	1:C:995:ARG:HH12	1.66	0.60
1:A:964:LYS:NZ	1:B:569:ILE:O	2.33	0.60
1:C:442:ASP:O	1:C:448:ASN:ND2	2.34	0.60
1:C:326:ILE:HD11	1:C:534:VAL:HB	1.84	0.60
1:C:345:THR:O	1:C:509:ARG:NH2	2.31	0.60
2:D:6:GLN:NE2	2:D:86:TYR:O	2.35	0.60
1:A:115:GLN:HE21	1:A:130:VAL:HG12	1.66	0.59
1:B:43:PHE:HB3	1:C:566:GLY:HA2	1.84	0.59
1:C:866:THR:HG22	1:C:869:MET:HE3	1.84	0.59
1:C:656:VAL:HG12	1:C:658:ASN:H	1.67	0.59
2:D:44:PRO:O	3:E:112:TRP:NE1	2.35	0.59
1:A:247:SER:OG	1:A:259:THR:OG1	2.20	0.59
1:A:96:GLU:OE2	1:A:100:ILE:N	2.30	0.59
1:C:409:GLN:HA	1:C:414:GLN:HG2	1.84	0.59
3:I:39:GLN:HB2	3:I:45:LEU:HD23	1.85	0.58
2:H:34:ALA:HB2	3:I:108:TYR:HD1	1.66	0.58
1:A:786:LYS:HE3	1:A:786:LYS:HA	1.85	0.58
1:B:43:PHE:HA	1:C:563:GLN:NE2	2.18	0.58
1:B:898:PHE:HZ	1:B:1050:MET:HE1	1.67	0.58
1:A:740:MET:HE3	1:B:319:ARG:HH21	1.68	0.58
1:A:869:MET:HG2	1:B:699:LEU:HD21	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:LEU:HB3	1:C:243:ALA:HA	1.86	0.58
1:C:1092:GLU:O	1:C:1107:ARG:NE	2.37	0.58
2:D:90:GLN:HE22	2:D:97:VAL:HG21	1.69	0.58
1:B:357:ARG:NH1	1:B:358:ILE:O	2.37	0.57
1:C:986:PRO:O	1:C:989:ALA:N	2.36	0.57
1:B:53:ASP:OD1	1:B:54:LEU:N	2.36	0.57
3:I:50:CYS:HA	3:I:71:ILE:HD13	1.86	0.57
1:A:873:TYR:CE1	1:B:699:LEU:HD12	2.39	0.57
1:B:197:ILE:HB	1:B:202:LYS:HZ1	1.70	0.57
1:A:1135:ASN:OD1	1:A:1136:THR:N	2.37	0.57
1:C:566:GLY:HA3	1:C:575:ALA:HB3	1.87	0.57
3:E:68:ARG:NH2	3:E:90:ASP:OD2	2.37	0.57
1:A:344:ALA:HB3	1:A:347:PHE:HE1	1.69	0.57
1:C:95:ILE:HD13	1:C:189:LEU:HD13	1.86	0.57
1:B:642:VAL:HG22	1:B:651:ILE:HG22	1.87	0.57
1:B:42:VAL:HG11	1:C:567:ARG:HH21	1.69	0.57
1:B:332:ILE:HD12	1:B:527:PRO:HA	1.86	0.57
3:G:50:CYS:HA	3:G:71:ILE:HD13	1.87	0.57
1:A:453:TYR:HB3	1:A:495:TYR:CE1	2.39	0.56
1:A:520:ALA:HB3	1:A:521:PRO:HD3	1.86	0.56
1:C:353:TRP:O	1:C:466:ARG:NH2	2.37	0.56
2:D:45:LYS:HA	3:E:112:TRP:HE1	1.70	0.56
3:E:47:TRP:HE1	3:E:50:CYS:HB2	1.71	0.56
2:F:43:PRO:HB3	3:G:112:TRP:HE3	1.70	0.56
1:A:490:PHE:HB3	1:A:493:ARG:HH22	1.69	0.56
1:B:36:VAL:HG23	1:B:222:ALA:HA	1.86	0.56
1:B:418:ILE:HA	1:B:422:ASN:HD22	1.70	0.56
1:B:790:LYS:NZ	1:C:702:GLU:OE2	2.39	0.56
1:C:898:PHE:HZ	1:C:1050:MET:HE1	1.71	0.56
3:G:35:CYS:HA	3:G:50:CYS:CB	2.36	0.56
1:A:280:ASN:OD1	1:A:284:THR:N	2.38	0.56
3:E:38:ARG:HB3	3:E:48:ILE:HD11	1.87	0.56
2:F:44:PRO:HD2	3:G:112:TRP:CZ3	2.40	0.56
1:B:732:THR:HB	1:B:955:ASN:ND2	2.21	0.56
3:I:34:MET:HB2	3:I:98:ARG:HH11	1.71	0.56
1:A:201:PHE:HB3	1:A:229:LEU:HB2	1.87	0.56
1:A:448:ASN:HB3	1:A:497:PHE:HD2	1.70	0.56
1:C:992:GLN:OE1	1:C:995:ARG:NH2	2.39	0.56
1:A:310:LYS:HG2	1:A:664:ILE:HD11	1.86	0.55
1:B:751:ASN:O	1:B:755:GLN:NE2	2.38	0.55
2:H:47:LEU:HD13	2:H:62:PHE:CD1	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:644:GLN:NE2	1:C:648:GLY:O	2.39	0.55
2:H:46:LEU:HD21	2:H:49:TYR:HB3	1.87	0.55
2:D:98:ASP:OD1	2:D:99:ASN:N	2.40	0.55
1:B:456:PHE:HD1	1:B:491:PRO:HA	1.72	0.55
1:C:115:GLN:NE2	1:C:165:ASN:O	2.40	0.55
2:D:5:THR:OG1	2:D:24:GLN:NE2	2.39	0.55
2:F:61:ARG:NH2	2:F:79:GLU:OE2	2.40	0.55
3:I:35:CYS:SG	3:I:47:TRP:NE1	2.80	0.55
3:I:28:PHE:HB3	3:I:53:THR:HB	1.88	0.55
1:A:130:VAL:HG21	1:A:231:ILE:HD12	1.88	0.55
2:H:90:GLN:NE2	2:H:93:SER:O	2.40	0.55
1:B:409:GLN:HA	1:B:414:GLN:HG2	1.89	0.55
3:G:109:PHE:HB3	3:G:112:TRP:HE1	1.72	0.55
1:C:356:LYS:HB3	1:C:397:ALA:HB3	1.89	0.54
3:G:38:ARG:HH22	3:G:40:ALA:HB2	1.71	0.54
3:I:109:PHE:HB3	3:I:112:TRP:HE1	1.71	0.54
1:B:216:LEU:HD12	1:B:217:PRO:HD2	1.88	0.54
1:A:401:VAL:HG22	1:A:509:ARG:HG2	1.89	0.54
3:E:33:ASP:HB3	3:E:99:ASP:HB3	1.90	0.54
1:B:128:ILE:HB	1:B:170:TYR:HB3	1.89	0.54
1:B:344:ALA:HB3	1:B:347:PHE:HE1	1.73	0.54
1:B:401:VAL:HG22	1:B:509:ARG:HG2	1.88	0.54
2:F:43:PRO:HB3	3:G:112:TRP:CE3	2.42	0.54
1:B:544:ASN:HD21	1:B:579:PRO:HB3	1.72	0.54
1:B:332:ILE:HG22	1:B:362:VAL:HG11	1.88	0.54
1:A:22:THR:O	1:A:78:ARG:NH1	2.40	0.54
1:A:358:ILE:HB	1:A:395:VAL:HB	1.90	0.54
1:B:821:LEU:HD11	1:B:939:SER:HB2	1.90	0.54
1:B:159:VAL:HG23	1:B:160:TYR:HD2	1.73	0.53
1:B:736:VAL:HG12	1:B:858:LEU:HD22	1.89	0.53
1:C:82:PRO:HG2	1:C:84:LEU:HD21	1.89	0.53
1:B:972:ALA:HB1	1:B:992:GLN:NE2	2.22	0.53
1:B:1135:ASN:OD1	1:B:1136:THR:N	2.41	0.53
1:C:732:THR:HB	1:C:955:ASN:ND2	2.22	0.53
1:C:906:PHE:CD2	1:C:916:LEU:HB2	2.43	0.53
1:B:106:PHE:HB2	1:B:117:LEU:HB3	1.90	0.53
2:H:59:PRO:HG2	2:H:62:PHE:CD2	2.44	0.53
1:B:22:THR:O	1:B:78:ARG:NH1	2.42	0.53
1:B:342:PHE:HZ	1:B:434:ILE:HD13	1.73	0.53
2:F:38:GLN:HB2	2:F:42:GLN:O	2.08	0.53
2:F:43:PRO:HG3	3:G:95:PHE:CE1	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:972:ALA:HB1	1:B:992:GLN:HE21	1.73	0.53
1:C:520:ALA:HB3	1:C:521:PRO:HD3	1.90	0.53
3:E:38:ARG:HH21	3:E:92:ALA:HB2	1.74	0.53
3:G:35:CYS:SG	3:G:47:TRP:NE1	2.81	0.53
1:A:391:CYS:HA	1:A:525:CYS:HA	1.91	0.52
1:A:898:PHE:HZ	1:A:1050:MET:HE1	1.73	0.52
1:C:142:ASP:OD2	1:C:246:ARG:NH2	2.42	0.52
2:H:37:GLN:HE21	2:H:84:ALA:HB3	1.73	0.52
1:A:418:ILE:HA	1:A:422:ASN:HD22	1.74	0.52
1:B:31:SER:N	1:B:60:SER:O	2.41	0.52
1:C:379:CYS:HB2	1:C:432:CYS:HA	1.91	0.52
1:A:318:PHE:HZ	1:A:615:VAL:HG11	1.73	0.52
1:B:346:ARG:HE	3:G:103:TYR:HB3	1.73	0.52
1:C:164:ASN:OD1	1:C:165:ASN:N	2.42	0.52
1:B:412:PRO:HG3	1:B:429:PHE:HB3	1.92	0.52
1:C:90:VAL:HG13	1:C:267:VAL:HG23	1.92	0.52
1:A:310:LYS:HG3	1:A:600:PRO:HA	1.92	0.52
2:H:6:GLN:NE2	2:H:86:TYR:O	2.42	0.52
1:A:1086:LYS:HD2	1:A:1122:VAL:HG11	1.90	0.52
1:B:722:VAL:HG22	1:B:1065:VAL:HG22	1.92	0.52
1:C:130:VAL:HG21	1:C:231:ILE:HD12	1.92	0.52
1:C:703:ASN:OD1	1:C:704:SER:N	2.43	0.52
1:B:395:VAL:HG13	1:B:515:PHE:HE1	1.75	0.52
1:A:454:ARG:HD3	1:A:457:ARG:HB2	1.92	0.52
1:C:280:ASN:OD1	1:C:284:THR:N	2.40	0.52
1:C:350:VAL:HG22	1:C:422:ASN:HB3	1.92	0.52
3:G:38:ARG:NH1	3:G:46:GLU:OE1	2.43	0.52
1:C:453:TYR:CE1	1:C:493:ARG:HB3	2.45	0.51
1:A:346:ARG:HB3	2:D:32:TRP:CZ2	2.45	0.51
1:B:546:LEU:HD21	1:B:573:THR:HG21	1.92	0.51
3:I:33:ASP:HB3	3:I:99:ASP:HB3	1.93	0.51
1:A:364:ASP:HB3	1:A:527:PRO:HB3	1.93	0.51
1:B:57:PRO:HG3	1:B:273:ARG:HD2	1.91	0.51
1:B:873:TYR:CE1	1:C:699:LEU:HG	2.46	0.51
1:C:401:VAL:HG22	1:C:509:ARG:HG2	1.93	0.51
1:C:985:ASP:OD1	1:C:986:PRO:HD2	2.10	0.51
3:I:48:ILE:HG13	3:I:49:ALA:N	2.26	0.51
1:B:102:ARG:HG3	1:B:141:LEU:HD12	1.93	0.51
2:D:91:THR:HG21	3:E:108:TYR:CE1	2.45	0.51
2:F:91:THR:HG21	3:G:108:TYR:CE1	2.46	0.51
3:G:32:TYR:HB2	3:G:98:ARG:HH21	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:331:ASN:O	1:C:333:THR:HG23	2.11	0.51
3:I:70:THR:N	3:I:82:GLN:O	2.40	0.51
1:B:94:SER:HA	1:B:265:TYR:HA	1.93	0.51
1:B:742:ILE:O	1:B:1000:ARG:NH1	2.43	0.51
1:C:493:ARG:HH21	1:C:495:TYR:HA	1.76	0.51
3:E:37:VAL:HG12	3:E:47:TRP:HA	1.93	0.50
2:F:90:GLN:HE21	2:F:100:VAL:HG12	1.75	0.50
1:A:231:ILE:HG22	1:A:233:ILE:HG23	1.94	0.50
1:B:86:PHE:HA	1:B:238:PHE:HE1	1.76	0.50
1:B:329:PHE:H	1:B:530:SER:HB2	1.75	0.50
1:C:468:ILE:HB	2:H:67:SER:HB3	1.93	0.50
1:C:763:LEU:HG	1:C:1008:VAL:HG21	1.92	0.50
1:B:234:ASN:HB2	5:B:1302:NAG:O5	2.12	0.50
1:A:915:VAL:O	1:A:919:ASN:ND2	2.41	0.50
1:C:80:ASP:O	1:C:265:TYR:OH	2.25	0.50
1:C:930:ALA:O	1:C:934:ILE:HG12	2.11	0.50
1:C:124:THR:OG1	1:C:125:ASN:OD1	2.24	0.50
1:C:139:PRO:HB3	1:C:159:VAL:HA	1.93	0.50
1:B:767:LEU:HD21	1:B:1008:VAL:HG22	1.94	0.50
1:A:821:LEU:HD11	1:A:939:SER:HB2	1.93	0.49
3:G:19:LEU:HD13	3:G:83:MET:HE3	1.94	0.49
1:B:377:PHE:CE2	1:B:384:PRO:HB3	2.47	0.49
2:F:49:TYR:HE1	2:F:55:ALA:HA	1.78	0.49
1:B:520:ALA:HB3	1:B:521:PRO:HD3	1.95	0.49
1:B:906:PHE:CD2	1:B:916:LEU:HB2	2.48	0.49
2:D:31:ASN:O	2:D:32:TRP:C	2.55	0.49
2:F:6:GLN:NE2	2:F:86:TYR:O	2.45	0.49
3:I:68:ARG:HH21	3:I:86:LEU:HA	1.78	0.49
3:G:109:PHE:CB	3:G:112:TRP:HE1	2.25	0.49
1:B:170:TYR:CZ	1:B:172:SER:HB2	2.48	0.49
1:B:357:ARG:NH1	1:B:395:VAL:H	2.10	0.49
1:B:462:LYS:HB2	1:B:465:GLU:HG3	1.94	0.49
1:B:974:SER:HB3	1:B:980:ILE:HD11	1.94	0.49
1:C:346:ARG:HE	3:I:103:TYR:HB3	1.77	0.49
3:I:64:TRP:CD1	3:I:64:TRP:H	2.30	0.49
1:A:132:GLU:OE1	1:A:165:ASN:ND2	2.46	0.49
1:C:922:LEU:HD11	1:C:926:GLN:HE21	1.77	0.49
1:C:642:VAL:HG22	1:C:651:ILE:HD12	1.94	0.48
1:C:190:ARG:HB3	1:C:192:PHE:HE1	1.77	0.48
1:A:1126:CYS:HB2	1:A:1132:ILE:HD13	1.94	0.48
1:B:377:PHE:HE2	1:B:384:PRO:HB3	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:49:TYR:HE1	2:H:55:ALA:HA	1.78	0.48
2:H:59:PRO:HG2	2:H:62:PHE:CE2	2.48	0.48
1:A:68:ILE:HG22	1:A:78:ARG:HB2	1.95	0.48
1:A:412:PRO:HG3	1:A:429:PHE:HB3	1.94	0.48
1:C:64:TRP:CD1	1:C:266:TYR:HE1	2.32	0.48
1:C:442:ASP:OD2	1:C:509:ARG:NE	2.45	0.48
1:A:740:MET:HE3	1:B:319:ARG:NH2	2.28	0.48
1:A:421:TYR:CD1	1:A:457:ARG:HB3	2.49	0.48
2:F:36:TYR:HA	2:F:45:LYS:O	2.14	0.48
3:G:68:ARG:HG3	3:G:69:PHE:HD1	1.79	0.48
3:E:98:ARG:HH12	3:E:111:LEU:HD13	1.79	0.48
1:A:431:GLY:HA2	1:A:515:PHE:CD2	2.49	0.48
1:B:383:SER:HB3	1:B:386:LYS:HG2	1.95	0.48
1:C:294:ASP:OD1	1:C:294:ASP:N	2.47	0.48
3:E:18:THR:HG23	3:E:82:GLN:HB3	1.95	0.48
3:G:51:ILE:HA	3:G:59:ILE:HA	1.95	0.48
1:B:328:ARG:NE	1:B:578:ASP:OD1	2.39	0.48
1:B:900:MET:HE1	1:C:1094:VAL:HG23	1.95	0.48
1:B:973:ILE:HG12	1:B:992:GLN:HE22	1.79	0.48
2:F:85:THR:HG22	2:F:106:GLU:HB3	1.96	0.48
3:I:88:ALA:HA	3:I:120:VAL:HG11	1.96	0.48
3:E:41:PRO:HD3	3:E:92:ALA:HA	1.95	0.47
1:A:995:ARG:NH2	1:C:994:ASP:OD2	2.46	0.47
2:D:94:SER:OG	3:E:106:PRO:HD2	2.14	0.47
3:G:32:TYR:O	3:G:53:THR:OG1	2.32	0.47
1:A:930:ALA:O	1:A:934:ILE:HG12	2.13	0.47
1:A:968:SER:HB3	1:A:970:PHE:CE2	2.49	0.47
1:B:64:TRP:CD1	1:B:266:TYR:HE1	2.33	0.47
1:B:989:ALA:O	1:B:993:ILE:HG12	2.14	0.47
1:A:462:LYS:HG2	1:A:465:GLU:OE1	2.14	0.47
1:C:24:LEU:HD12	1:C:25:PRO:HD2	1.96	0.47
1:A:29:THR:OG1	1:A:215:ASP:OD1	2.28	0.47
1:A:247:SER:N	1:A:257:GLY:O	2.44	0.47
1:A:658:ASN:ND2	1:A:660:TYR:OH	2.48	0.47
1:B:68:ILE:HG22	1:B:78:ARG:HB2	1.96	0.47
1:B:403:ARG:HH22	1:B:505:HIS:CE1	2.32	0.47
1:C:125:ASN:HB2	1:C:172:SER:O	2.14	0.47
1:C:750:SER:O	1:C:754:LEU:HG	2.15	0.47
3:E:98:ARG:HH22	3:E:111:LEU:HD13	1.80	0.47
3:I:34:MET:HB2	3:I:98:ARG:NH1	2.29	0.47
1:B:96:GLU:OE2	1:B:101:ILE:N	2.33	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:334:ASN:O	1:C:362:VAL:HG12	2.15	0.47
1:C:740:MET:HE2	1:C:740:MET:HA	1.95	0.47
1:B:246:ARG:HD3	1:B:258:TRP:HB3	1.95	0.47
1:B:334:ASN:HB3	1:B:361:CYS:HA	1.97	0.47
2:H:94:SER:OG	3:I:106:PRO:HD2	2.15	0.47
3:I:35:CYS:HA	3:I:50:CYS:HB2	1.97	0.47
1:A:246:ARG:HA	1:A:258:TRP:HA	1.97	0.47
1:B:129:LYS:HE3	1:B:133:PHE:HZ	1.80	0.47
1:C:986:PRO:HA	1:C:989:ALA:HB3	1.96	0.47
2:D:61:ARG:NH2	2:D:79:GLU:OE2	2.48	0.47
1:C:125:ASN:ND2	1:C:171:VAL:HG13	2.30	0.46
3:E:8:GLY:HA2	3:E:118:VAL:HG22	1.97	0.46
1:A:562:PHE:HB2	1:C:41:LYS:NZ	2.31	0.46
1:A:566:GLY:N	1:A:575:ALA:O	2.49	0.46
1:C:335:LEU:HD13	1:C:362:VAL:HG13	1.98	0.46
3:E:32:TYR:O	3:E:53:THR:OG1	2.33	0.46
1:C:281:GLU:H	1:C:281:GLU:CD	2.23	0.46
1:C:732:THR:HB	1:C:955:ASN:HD22	1.81	0.46
1:A:444:LYS:HA	1:A:444:LYS:HD3	1.68	0.46
1:A:763:LEU:HD22	1:A:1008:VAL:HG21	1.98	0.46
1:B:226:LEU:HG	1:B:227:VAL:HG23	1.98	0.46
2:D:20:THR:HG23	2:D:22:LYS:HZ1	1.80	0.46
2:F:34:ALA:HB3	2:F:89:GLN:HB3	1.97	0.46
3:E:63:SER:HA	3:E:66:LYS:HE3	1.98	0.46
3:G:48:ILE:HG12	3:G:64:TRP:CH2	2.51	0.46
1:A:81:ASN:O	1:A:239:GLN:NE2	2.43	0.46
1:A:442:ASP:OD1	1:A:451:TYR:OH	2.20	0.46
1:B:106:PHE:HD2	1:B:117:LEU:HD22	1.81	0.46
1:B:344:ALA:HB3	1:B:347:PHE:CE1	2.51	0.46
1:C:128:ILE:HG21	1:C:229:LEU:HD21	1.98	0.46
2:H:38:GLN:NE2	3:I:39:GLN:OE1	2.45	0.46
3:I:48:ILE:HG13	3:I:49:ALA:H	1.80	0.46
3:G:63:SER:O	3:G:66:LYS:NZ	2.37	0.46
2:H:22:LYS:HA	2:H:72:THR:HA	1.98	0.46
1:A:126:VAL:HB	1:A:172:SER:O	2.16	0.46
1:C:132:GLU:HB3	1:C:164:ASN:OD1	2.16	0.46
1:B:523:THR:HG23	1:B:524:VAL:HG23	1.98	0.45
1:B:559:PHE:HB3	1:B:563:GLN:HB2	1.96	0.45
3:G:50:CYS:HA	3:G:71:ILE:HG21	1.97	0.45
1:A:457:ARG:NH1	1:A:467:ASP:OD2	2.50	0.45
1:C:1126:CYS:HB2	1:C:1132:ILE:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:34:MET:HG2	3:G:79:VAL:HG11	1.98	0.45
2:H:38:GLN:HB3	2:H:44:PRO:HB3	1.98	0.45
2:D:1:ASP:N	2:D:96:ASP:O	2.50	0.45
1:A:64:TRP:CD1	1:A:266:TYR:HE1	2.35	0.45
1:A:97:LYS:HG2	1:A:186:PHE:HD1	1.82	0.45
1:A:599:THR:HG22	1:A:608:VAL:HG12	1.98	0.45
1:A:699:LEU:HD22	1:C:873:TYR:CZ	2.51	0.45
1:B:865:LEU:HD13	1:B:869:MET:HE3	1.99	0.45
1:C:418:ILE:HA	1:C:422:ASN:HD22	1.81	0.45
2:D:59:PRO:HG2	2:D:62:PHE:HE1	1.80	0.45
3:G:38:ARG:HH21	3:G:92:ALA:HB2	1.82	0.45
2:H:63:LYS:HD3	2:H:74:THR:HB	1.99	0.45
1:A:866:THR:H	1:A:869:MET:HE3	1.82	0.45
1:A:902:MET:HB3	1:A:916:LEU:HD11	1.97	0.45
1:B:334:ASN:C	1:B:362:VAL:HG12	2.41	0.45
1:B:644:GLN:NE2	1:B:645:THR:O	2.45	0.45
1:C:231:ILE:HG22	1:C:233:ILE:HG23	1.99	0.45
1:A:388:ASN:O	1:A:526:GLY:HA3	2.16	0.45
1:C:129:LYS:HE3	1:C:133:PHE:HZ	1.80	0.45
1:C:187:LYS:O	1:C:210:ILE:N	2.43	0.45
1:A:311:GLY:HA2	1:A:664:ILE:HD12	1.99	0.45
1:A:345:THR:O	1:A:509:ARG:NH2	2.41	0.45
1:B:605:SER:OG	1:B:606:ASN:N	2.49	0.45
2:H:61:ARG:HD2	2:H:62:PHE:N	2.31	0.45
2:H:20:THR:HG22	2:H:74:THR:HA	1.98	0.45
1:A:699:LEU:HB3	1:C:873:TYR:HE1	1.81	0.45
1:B:712:ILE:O	1:B:1075:PHE:N	2.44	0.45
3:G:37:VAL:HG23	3:G:95:PHE:HB2	1.98	0.45
1:C:954:HIS:HB3	1:C:1014:ARG:CZ	2.47	0.45
1:B:142:ASP:CG	1:B:246:ARG:HE	2.24	0.44
1:A:53:ASP:HB3	1:A:55:PHE:CE2	2.52	0.44
1:A:902:MET:HE3	1:A:1049:LEU:HD13	2.00	0.44
1:C:125:ASN:HA	1:C:174:PRO:HD3	1.98	0.44
1:A:41:LYS:NZ	1:B:562:PHE:HB2	2.33	0.44
1:C:79:PHE:HE2	1:C:258:TRP:HE1	1.66	0.44
1:C:1082:CYS:HB2	1:C:1126:CYS:HB2	1.97	0.44
3:G:35:CYS:HA	3:G:50:CYS:HB2	1.97	0.44
1:B:890:ALA:HB1	1:C:1046:GLY:HA2	1.98	0.44
3:I:30:SER:HA	3:I:54:GLY:HA2	1.98	0.44
1:A:822:LEU:HD22	1:A:945:LEU:HD21	1.99	0.44
1:B:993:ILE:O	1:B:997:ILE:HG22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:MET:SD	1:A:153:MET:N	2.91	0.44
1:A:605:SER:OG	1:A:606:ASN:N	2.50	0.44
1:A:989:ALA:O	1:A:993:ILE:HG12	2.17	0.44
1:C:1086:LYS:HD2	1:C:1122:VAL:HG11	2.00	0.44
1:A:384:PRO:HA	1:A:387:LEU:HG	1.99	0.44
1:B:726:ILE:HG13	1:B:1061:VAL:HG22	1.99	0.44
1:B:409:GLN:OE1	1:B:418:ILE:N	2.51	0.44
1:B:457:ARG:HH12	1:B:461:LEU:HD21	1.81	0.44
1:C:540:ASN:HB3	1:C:547:LYS:NZ	2.33	0.44
1:A:382:VAL:HA	1:C:983:ARG:O	2.17	0.44
1:C:68:ILE:HG22	1:C:78:ARG:HB2	2.00	0.44
1:C:777:ASN:O	1:C:781:VAL:HG23	2.18	0.44
1:A:858:LEU:HD13	1:A:959:LEU:HD11	1.98	0.43
1:B:770:ILE:HD11	1:B:1012:LEU:HA	2.00	0.43
1:B:1102:TRP:HB2	1:B:1135:ASN:ND2	2.33	0.43
1:C:902:MET:HE2	1:C:902:MET:HA	2.00	0.43
2:F:2:ILE:HG13	2:F:27:GLU:HB2	2.00	0.43
1:A:451:TYR:HE1	2:D:32:TRP:CZ2	2.36	0.43
2:D:24:GLN:OE1	2:D:24:GLN:N	2.51	0.43
1:A:387:LEU:HA	1:A:390:LEU:HD12	1.99	0.43
1:A:900:MET:HE1	1:B:1077:THR:CG2	2.48	0.43
1:B:48:LEU:HD23	1:B:276:LEU:HD21	2.00	0.43
1:B:125:ASN:HD22	1:B:171:VAL:HG13	1.82	0.43
1:B:190:ARG:HB3	1:B:192:PHE:CE2	2.53	0.43
3:E:38:ARG:O	3:E:45:LEU:HD12	2.18	0.43
3:I:3:LEU:HD11	3:I:98:ARG:HB2	2.00	0.43
1:A:112:SER:HA	1:A:132:GLU:HB3	2.00	0.43
1:A:815:ARG:HB3	1:A:819:GLU:HG3	2.00	0.43
1:B:53:ASP:HB3	1:B:55:PHE:CE2	2.53	0.43
1:C:176:LEU:HG	1:C:207:HIS:HE1	1.83	0.43
1:B:106:PHE:HB3	1:B:235:ILE:HD12	2.00	0.43
1:B:804:GLN:NE2	1:B:935:GLN:OE1	2.35	0.43
1:B:1097:SER:HB3	1:B:1102:TRP:CD2	2.54	0.43
2:D:33:LEU:HD22	2:D:71:PHE:CD2	2.54	0.43
3:E:10:LEU:HD11	3:E:121:SER:HB3	2.01	0.43
3:I:41:PRO:HD3	3:I:92:ALA:HA	2.01	0.43
1:B:31:SER:HB3	1:B:60:SER:N	2.34	0.43
1:B:418:ILE:HA	1:B:422:ASN:ND2	2.34	0.43
1:C:131:CYS:HA	1:C:166:CYS:HA	2.00	0.43
2:H:33:LEU:HD13	2:H:71:PHE:CG	2.53	0.43
1:A:770:ILE:O	1:A:774:GLN:HG2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:1303:NAG:H82	1:C:558:LYS:HE3	2.00	0.43
3:G:32:TYR:HD1	3:G:99:ASP:O	2.02	0.43
1:A:43:PHE:CE1	1:A:283:GLY:HA3	2.54	0.43
1:A:393:THR:OG1	1:A:516:GLU:OE1	2.35	0.43
1:B:1091:ARG:HG3	1:B:1119:ASN:C	2.43	0.43
1:A:557:LYS:NZ	1:A:574:ASP:OD2	2.51	0.42
1:A:571:ASP:HB2	1:C:967:SER:HB2	2.01	0.42
1:A:1094:VAL:HG23	1:C:900:MET:HE1	1.99	0.42
2:H:110:LYS:HA	2:H:110:LYS:HD3	1.84	0.42
1:C:559:PHE:CE1	1:C:584:ILE:HG23	2.54	0.42
1:C:618:THR:OG1	1:C:619:GLU:OE1	2.29	0.42
1:A:142:ASP:OD2	1:A:246:ARG:NH2	2.32	0.42
1:A:659:SER:HB3	1:A:698:SER:HB3	2.00	0.42
1:B:280:ASN:ND2	1:B:282:ASN:OD1	2.52	0.42
1:B:444:LYS:HD3	1:B:444:LYS:HA	1.68	0.42
1:B:969:LYS:HZ2	1:B:974:SER:HA	1.83	0.42
1:C:403:ARG:HH22	1:C:505:HIS:CD2	2.37	0.42
1:C:344:ALA:HB3	1:C:347:PHE:HE1	1.83	0.42
1:C:444:LYS:HA	1:C:444:LYS:HD3	1.71	0.42
3:E:103:TYR:HB2	3:E:107:ASP:OD2	2.19	0.42
3:I:21:CYS:HB3	3:I:79:VAL:HG13	2.00	0.42
1:A:36:VAL:HG11	1:A:220:PHE:CZ	2.54	0.42
1:A:712:ILE:HG13	1:C:896:ILE:HG12	2.02	0.42
1:B:560:LEU:HB2	1:B:563:GLN:HG3	2.02	0.42
1:C:813:SER:OG	1:C:815:ARG:NH1	2.53	0.42
2:H:35:TRP:CD1	2:H:48:ILE:HB	2.54	0.42
1:A:125:ASN:CG	1:A:126:VAL:H	2.28	0.42
1:C:375:PHE:CE2	1:C:377:PHE:HB2	2.55	0.42
3:G:3:LEU:HA	3:G:3:LEU:HD23	1.82	0.42
1:A:457:ARG:HH12	1:A:461:LEU:HD21	1.85	0.42
1:A:1082:CYS:HB2	1:A:1126:CYS:HB2	1.87	0.42
1:B:15:CYS:N	1:B:158:ARG:HD3	2.34	0.42
2:F:38:GLN:O	2:F:84:ALA:HB1	2.20	0.42
3:I:37:VAL:HG23	3:I:95:PHE:HB2	2.01	0.42
1:A:1086:LYS:HB2	1:A:1086:LYS:HE2	1.90	0.42
1:B:92:PHE:O	1:B:192:PHE:N	2.36	0.42
1:C:1037:SER:OG	1:C:1043:CYS:SG	2.66	0.42
3:E:109:PHE:HB2	3:E:112:TRP:CH2	2.54	0.42
1:A:287:ASP:OD1	1:A:288:ALA:N	2.53	0.42
1:B:348:ALA:HB2	1:B:354:ASN:ND2	2.35	0.42
1:C:220:PHE:HE2	1:C:285:ILE:HG22	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:361:CYS:O	1:C:524:VAL:HA	2.18	0.42
1:C:1102:TRP:HB2	1:C:1135:ASN:ND2	2.35	0.42
1:A:24:LEU:HB2	1:A:78:ARG:CZ	2.50	0.42
1:A:68:ILE:HG23	1:A:78:ARG:H	1.83	0.42
1:A:1030:SER:OG	1:B:1041:ASP:HB2	2.19	0.42
1:A:82:PRO:HG2	1:A:84:LEU:HD21	2.01	0.41
1:A:580:GLN:C	1:A:582:LEU:H	2.28	0.41
1:B:41:LYS:O	1:C:564:GLN:N	2.52	0.41
1:B:896:ILE:HG12	1:C:712:ILE:HG13	2.02	0.41
3:I:36:TRP:H	3:I:50:CYS:HB3	1.85	0.41
3:I:68:ARG:NH2	3:I:86:LEU:HA	2.34	0.41
1:A:906:PHE:CE2	1:A:916:LEU:HD12	2.55	0.41
1:C:190:ARG:HB3	1:C:192:PHE:CE1	2.55	0.41
1:C:328:ARG:NH2	1:C:581:THR:OG1	2.54	0.41
3:I:35:CYS:HA	3:I:50:CYS:CB	2.50	0.41
1:C:316:SER:OG	1:C:317:ASN:N	2.53	0.41
1:C:419:ALA:HA	1:C:423:TYR:O	2.20	0.41
1:C:722:VAL:O	1:C:934:ILE:HD11	2.21	0.41
3:G:39:GLN:HB2	3:G:45:LEU:HD13	2.01	0.41
1:A:277:LEU:HD12	1:A:285:ILE:HD13	2.02	0.41
1:A:456:PHE:HB3	1:A:473:TYR:CG	2.55	0.41
1:B:433:VAL:HG12	1:B:512:VAL:HG13	2.01	0.41
1:B:748:GLU:OE1	1:B:748:GLU:N	2.37	0.41
1:B:759:PHE:O	1:B:763:LEU:HG	2.21	0.41
1:C:216:LEU:HG	1:C:266:TYR:CE2	2.56	0.41
1:C:383:SER:HB3	1:C:386:LYS:HG2	2.02	0.41
1:A:309:GLU:O	1:A:313:TYR:OH	2.32	0.41
1:A:418:ILE:HA	1:A:422:ASN:ND2	2.35	0.41
1:A:546:LEU:HD11	1:A:573:THR:HG21	2.03	0.41
1:A:900:MET:HE1	1:B:1077:THR:HG23	2.02	0.41
1:B:323:THR:OG1	1:B:537:LYS:NZ	2.43	0.41
1:C:30:ASN:HB3	1:C:32:PHE:CE1	2.56	0.41
1:C:393:THR:HG21	1:C:521:PRO:HG2	2.02	0.41
1:C:533:LEU:HD23	1:C:533:LEU:H	1.84	0.41
3:G:47:TRP:HH2	3:G:60:TYR:HB3	1.86	0.41
2:H:85:THR:HG22	2:H:106:GLU:HB3	2.02	0.41
1:A:954:HIS:HB3	1:A:1014:ARG:CZ	2.50	0.41
1:B:141:LEU:HD22	1:B:157:PHE:CD1	2.55	0.41
2:D:44:PRO:O	3:E:112:TRP:CD1	2.74	0.41
3:E:40:ALA:HB3	3:E:43:LYS:HB2	2.03	0.41
3:G:58:ASN:ND2	3:G:104:ALA:HA	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:VAL:HG12	1:A:512:VAL:HG13	2.02	0.41
1:A:742:ILE:HG23	1:A:743:CYS:N	2.36	0.41
2:D:30:SER:HB3	2:D:32:TRP:HD1	1.85	0.41
3:G:38:ARG:HD2	3:G:48:ILE:HD11	2.03	0.41
1:A:329:PHE:O	1:A:580:GLN:NE2	2.39	0.41
1:A:549:THR:HB	1:C:745:ASP:HB3	2.03	0.41
1:A:713:ALA:HB3	1:C:894:LEU:HB3	2.02	0.41
1:B:42:VAL:HG11	1:C:567:ARG:NH2	2.35	0.41
1:C:125:ASN:HD22	1:C:171:VAL:HG13	1.86	0.41
1:C:388:ASN:O	1:C:526:GLY:HA3	2.21	0.41
2:D:38:GLN:O	2:D:84:ALA:HB1	2.20	0.41
2:H:61:ARG:NH1	2:H:75:ILE:HG12	2.35	0.41
2:H:85:THR:HA	2:H:106:GLU:HA	2.03	0.41
1:A:15:CYS:N	1:A:158:ARG:HD3	2.36	0.41
1:A:107:GLY:H	1:A:235:ILE:HG23	1.86	0.41
1:A:350:VAL:HG22	1:A:422:ASN:HB3	2.03	0.41
1:A:763:LEU:HD21	1:A:1005:GLN:OE1	2.21	0.41
1:B:132:GLU:OE1	1:B:165:ASN:ND2	2.54	0.41
1:B:578:ASP:OD2	1:B:581:THR:N	2.49	0.41
1:C:170:TYR:CE2	1:C:172:SER:HB2	2.55	0.41
1:C:741:TYR:CE1	1:C:966:LEU:HD13	2.56	0.41
3:E:38:ARG:HD3	3:E:64:TRP:HH2	1.86	0.41
3:E:87:THR:O	3:E:120:VAL:HG11	2.21	0.41
2:F:38:GLN:OE1	2:F:44:PRO:HD3	2.21	0.41
1:A:598:ILE:HG23	1:A:664:ILE:HG21	2.03	0.40
1:A:392:PHE:CD1	1:A:515:PHE:HB3	2.56	0.40
1:A:699:LEU:HD21	1:C:869:MET:HB3	2.04	0.40
1:A:1031:GLU:OE2	1:B:1039:ARG:NE	2.54	0.40
1:B:44:ARG:NH1	1:B:279:TYR:OH	2.53	0.40
1:C:128:ILE:HD13	1:C:170:TYR:HD2	1.86	0.40
2:D:36:TYR:CE2	3:E:112:TRP:HH2	2.39	0.40
3:G:3:LEU:HD21	3:G:26:PHE:HZ	1.85	0.40
1:A:455:LEU:HD22	1:A:493:ARG:HG3	2.03	0.40
1:A:1074:ASN:HD22	4:V:1:NAG:C7	2.35	0.40
1:B:375:PHE:HA	1:B:436:TRP:HB3	2.03	0.40
1:C:117:LEU:HD12	1:C:231:ILE:HG21	2.03	0.40
2:D:15:VAL:HA	2:D:78:VAL:HG23	2.02	0.40
2:D:18:THR:HG22	2:D:76:ASN:O	2.20	0.40
2:D:37:GLN:OE1	2:D:47:LEU:HD11	2.22	0.40
1:A:281:GLU:OE2	5:A:1303:NAG:O6	2.31	0.40
2:D:6:GLN:HE22	2:D:86:TYR:C	2.29	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:5:GLU:HG3	3:G:115:GLY:HA2	2.04	0.40
1:A:722:VAL:HG22	1:A:1065:VAL:HG22	2.02	0.40
1:B:83:VAL:HG11	1:B:237:ARG:HE	1.86	0.40
1:B:322:PRO:HG3	1:B:549:THR:HG21	2.03	0.40
1:B:364:ASP:OD2	1:B:366:SER:OG	2.38	0.40
1:B:392:PHE:CE1	1:B:517:LEU:HD13	2.56	0.40
1:B:421:TYR:HB3	1:B:454:ARG:HG2	2.03	0.40
1:B:996:LEU:HD23	1:B:996:LEU:HA	1.92	0.40
2:D:29:ILE:HD11	2:D:33:LEU:HD13	2.03	0.40
2:D:79:GLU:OE1	2:D:79:GLU:N	2.55	0.40
3:E:15:ALA:O	3:E:85:SER:N	2.54	0.40
2:F:38:GLN:HB3	2:F:44:PRO:CA	2.50	0.40
2:F:110:LYS:HA	2:F:110:LYS:HD3	1.84	0.40
3:I:83:MET:SD	3:I:83:MET:N	2.94	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	1016/1286 (79%)	982 (97%)	33 (3%)	1 (0%)	48 78
1	B	1017/1286 (79%)	977 (96%)	39 (4%)	1 (0%)	48 78
1	C	1017/1286 (79%)	973 (96%)	43 (4%)	1 (0%)	48 78
2	D	109/111 (98%)	104 (95%)	4 (4%)	1 (1%)	14 47
2	F	109/111 (98%)	102 (94%)	7 (6%)	0	100 100
2	H	109/111 (98%)	103 (94%)	6 (6%)	0	100 100
3	E	120/122 (98%)	116 (97%)	4 (3%)	0	100 100
3	G	120/122 (98%)	117 (98%)	3 (2%)	0	100 100
3	I	120/122 (98%)	116 (97%)	4 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	3737/4557 (82%)	3590 (96%)	143 (4%)	4 (0%)	50 78

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	520	ALA
1	B	520	ALA
1	C	520	ALA
2	D	32	TRP

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	908/1116 (81%)	908 (100%)	0	100 100
1	B	907/1116 (81%)	907 (100%)	0	100 100
1	C	906/1116 (81%)	906 (100%)	0	100 100
2	D	90/90 (100%)	90 (100%)	0	100 100
2	F	90/90 (100%)	90 (100%)	0	100 100
2	H	90/90 (100%)	90 (100%)	0	100 100
3	E	96/96 (100%)	96 (100%)	0	100 100
3	G	96/96 (100%)	96 (100%)	0	100 100
3	I	96/96 (100%)	96 (100%)	0	100 100
All	All	3279/3906 (84%)	3279 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	99	ASN

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Mol	Chain	Res	Type
1	A	164	ASN
1	A	165	ASN
1	A	901	GLN
1	A	926	GLN
1	A	965	GLN
1	A	992	GLN
1	A	1083	HIS
1	B	23	GLN
1	B	115	GLN
1	B	165	ASN
1	B	321	GLN
1	B	360	ASN
1	B	414	GLN
1	B	564	GLN
1	B	613	GLN
1	B	914	ASN
1	B	1010	GLN
1	B	1106	GLN
1	C	23	GLN
1	C	115	GLN
1	C	919	ASN
1	C	928	ASN
2	D	38	GLN
2	D	90	GLN
2	F	24	GLN
2	H	70	GLN
2	H	90	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

30 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	$\#\|Z\| > 2$
4	NAG	J	1	1,4	14,14,15	0.29	0	17,19,21	0.51	0
4	NAG	J	2	4	14,14,15	0.17	0	17,19,21	0.46	0
4	NAG	K	1	1,4	14,14,15	0.24	0	17,19,21	0.52	0
4	NAG	K	2	4	14,14,15	0.26	0	17,19,21	0.37	0
4	NAG	L	1	1,4	14,14,15	0.20	0	17,19,21	0.48	0
4	NAG	L	2	4	14,14,15	0.23	0	17,19,21	0.38	0
4	NAG	M	1	1,4	14,14,15	0.20	0	17,19,21	0.45	0
4	NAG	M	2	4	14,14,15	0.24	0	17,19,21	0.43	0
4	NAG	N	1	1,4	14,14,15	0.19	0	17,19,21	0.45	0
4	NAG	N	2	4	14,14,15	0.22	0	17,19,21	0.42	0
4	NAG	O	1	1,4	14,14,15	0.25	0	17,19,21	0.43	0
4	NAG	O	2	4	14,14,15	0.19	0	17,19,21	0.44	0
4	NAG	P	1	1,4	14,14,15	0.25	0	17,19,21	0.58	0
4	NAG	P	2	4	14,14,15	0.20	0	17,19,21	0.45	0
4	NAG	Q	1	1,4	14,14,15	0.31	0	17,19,21	0.63	1 (5%)
4	NAG	Q	2	4	14,14,15	0.26	0	17,19,21	0.43	0
4	NAG	R	1	1,4	14,14,15	0.19	0	17,19,21	0.47	0
4	NAG	R	2	4	14,14,15	0.19	0	17,19,21	0.42	0
4	NAG	S	1	1,4	14,14,15	0.22	0	17,19,21	0.46	0
4	NAG	S	2	4	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	T	1	1,4	14,14,15	0.18	0	17,19,21	0.46	0
4	NAG	T	2	4	14,14,15	0.21	0	17,19,21	0.42	0
4	NAG	U	1	1,4	14,14,15	0.21	0	17,19,21	0.56	0
4	NAG	U	2	4	14,14,15	0.21	0	17,19,21	0.40	0
4	NAG	V	1	1,4	14,14,15	0.21	0	17,19,21	0.38	0
4	NAG	V	2	4	14,14,15	0.19	0	17,19,21	0.50	0
4	NAG	W	1	1,4	14,14,15	0.19	0	17,19,21	0.45	0
4	NAG	W	2	4	14,14,15	0.21	0	17,19,21	0.42	0
4	NAG	X	1	1,4	14,14,15	0.20	0	17,19,21	0.42	0
4	NAG	X	2	4	14,14,15	0.20	0	17,19,21	0.45	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	J	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	J	2	4	-	2/6/23/26	0/1/1/1
4	NAG	K	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	K	2	4	-	1/6/23/26	0/1/1/1
4	NAG	L	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	L	2	4	-	2/6/23/26	0/1/1/1
4	NAG	M	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	M	2	4	-	2/6/23/26	0/1/1/1
4	NAG	N	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	N	2	4	-	0/6/23/26	0/1/1/1
4	NAG	O	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	O	2	4	-	2/6/23/26	0/1/1/1
4	NAG	P	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	P	2	4	-	2/6/23/26	0/1/1/1
4	NAG	Q	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	Q	2	4	-	2/6/23/26	0/1/1/1
4	NAG	R	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	R	2	4	-	2/6/23/26	0/1/1/1
4	NAG	S	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	S	2	4	-	0/6/23/26	0/1/1/1
4	NAG	T	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	T	2	4	-	0/6/23/26	0/1/1/1
4	NAG	U	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	U	2	4	-	2/6/23/26	0/1/1/1
4	NAG	V	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	V	2	4	-	2/6/23/26	0/1/1/1
4	NAG	W	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	W	2	4	-	0/6/23/26	0/1/1/1
4	NAG	X	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	X	2	4	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	Q	1	NAG	C1-O5-C5	2.09	115.02	112.19

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	J	1	NAG	O5-C5-C6-O6
4	Q	1	NAG	O5-C5-C6-O6
4	L	2	NAG	O5-C5-C6-O6
4	K	1	NAG	O5-C5-C6-O6
4	M	2	NAG	O5-C5-C6-O6
4	P	1	NAG	O5-C5-C6-O6
4	Q	2	NAG	C4-C5-C6-O6
4	J	1	NAG	C4-C5-C6-O6
4	Q	1	NAG	C4-C5-C6-O6
4	O	1	NAG	O5-C5-C6-O6
4	M	1	NAG	O5-C5-C6-O6
4	S	1	NAG	O5-C5-C6-O6
4	Q	2	NAG	O5-C5-C6-O6
4	P	2	NAG	O5-C5-C6-O6
4	N	1	NAG	O5-C5-C6-O6
4	O	1	NAG	C4-C5-C6-O6
4	L	2	NAG	C4-C5-C6-O6
4	P	2	NAG	C4-C5-C6-O6
4	X	1	NAG	O5-C5-C6-O6
4	M	1	NAG	C4-C5-C6-O6
4	S	1	NAG	C4-C5-C6-O6
4	K	1	NAG	C4-C5-C6-O6
4	M	2	NAG	C4-C5-C6-O6
4	P	1	NAG	C4-C5-C6-O6
4	X	1	NAG	C4-C5-C6-O6
4	T	1	NAG	O5-C5-C6-O6
4	U	2	NAG	O5-C5-C6-O6
4	N	1	NAG	C4-C5-C6-O6
4	T	1	NAG	C4-C5-C6-O6
4	U	2	NAG	C4-C5-C6-O6
4	O	2	NAG	C4-C5-C6-O6
4	O	2	NAG	O5-C5-C6-O6
4	K	2	NAG	O5-C5-C6-O6
4	V	1	NAG	O5-C5-C6-O6
4	R	2	NAG	C4-C5-C6-O6
4	V	2	NAG	C4-C5-C6-O6
4	R	2	NAG	O5-C5-C6-O6
4	V	2	NAG	O5-C5-C6-O6
4	J	2	NAG	C4-C5-C6-O6
4	J	2	NAG	O5-C5-C6-O6
4	J	1	NAG	C3-C2-N2-C7
4	P	1	NAG	C3-C2-N2-C7

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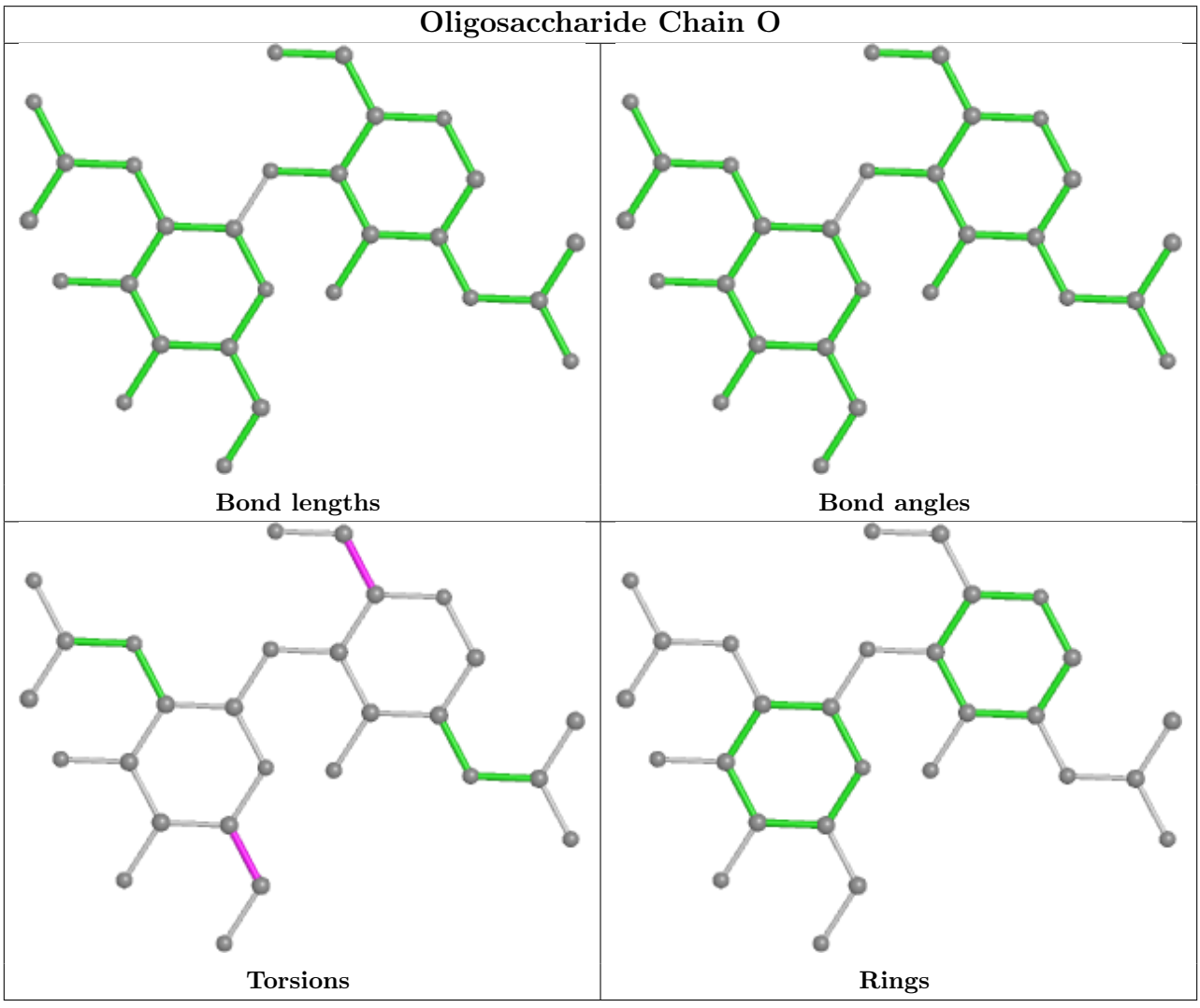
Mol	Chain	Res	Type	Atoms
4	U	1	NAG	C3-C2-N2-C7
4	X	2	NAG	C4-C5-C6-O6

There are no ring outliers.

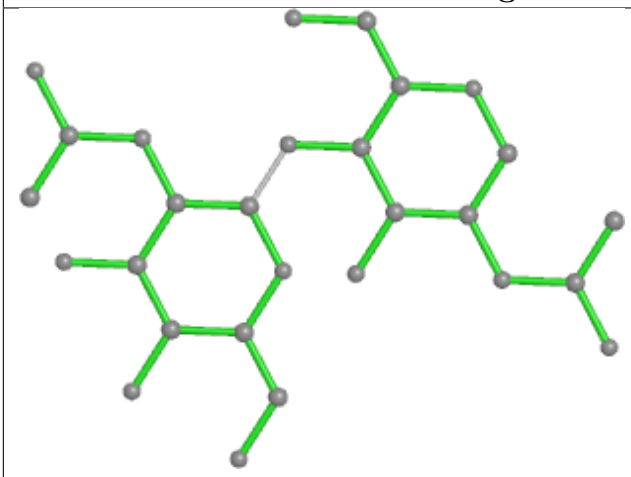
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	V	1	NAG	1	0

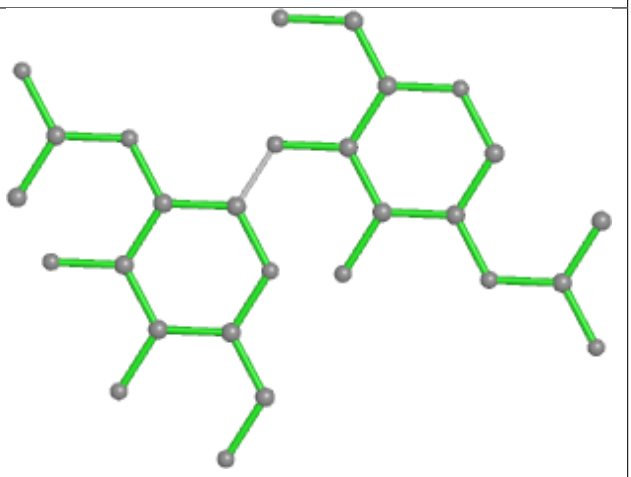
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



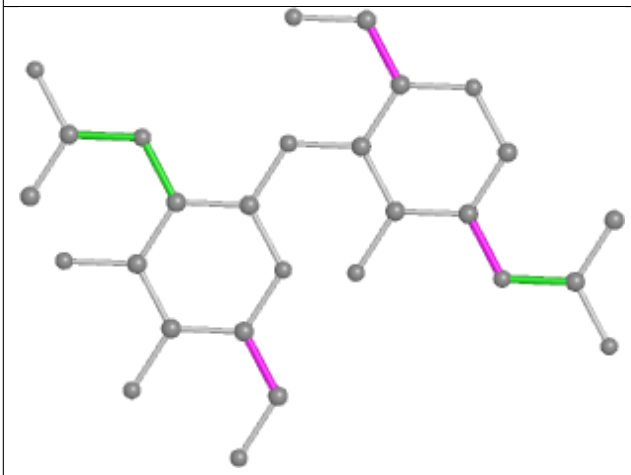
Oligosaccharide Chain J



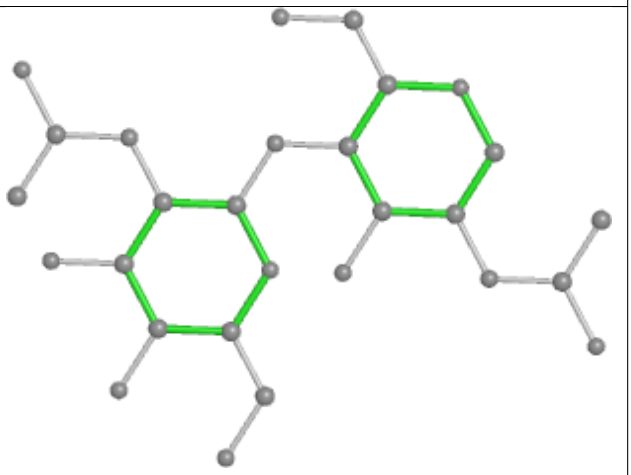
Bond lengths



Bond angles

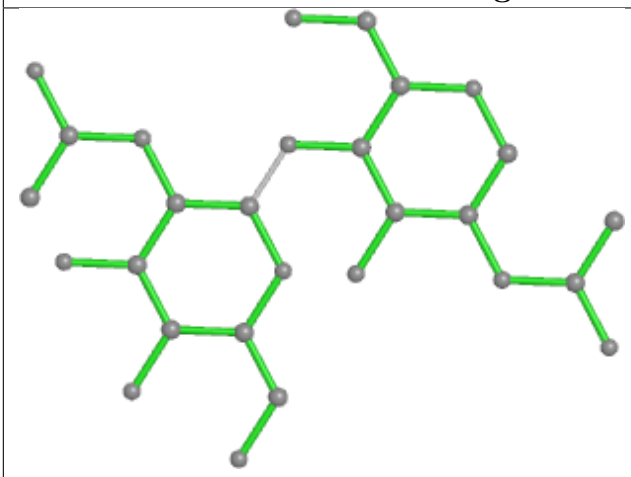


Torsions

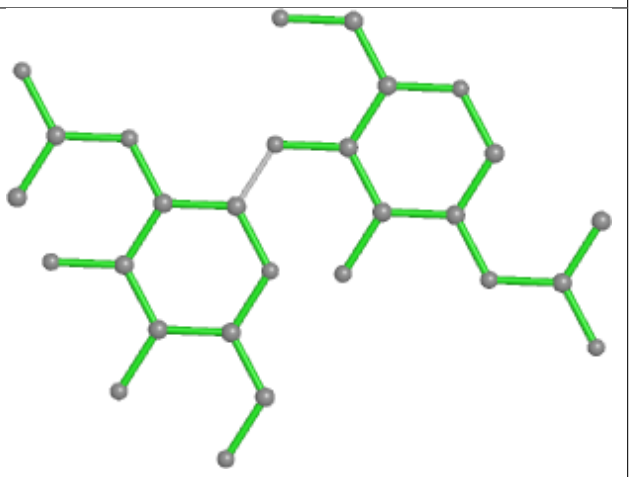


Rings

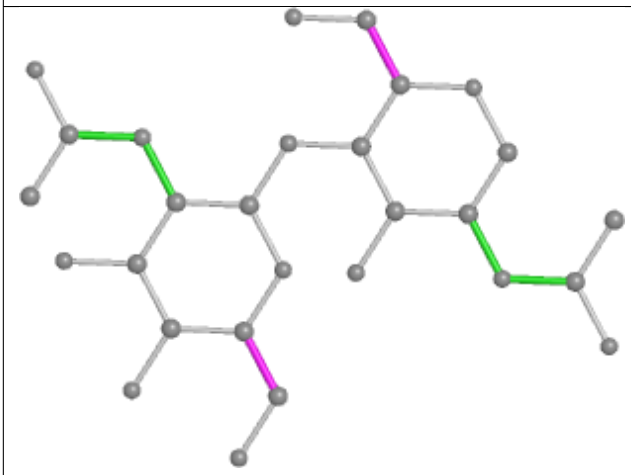
Oligosaccharide Chain K



Bond lengths



Bond angles



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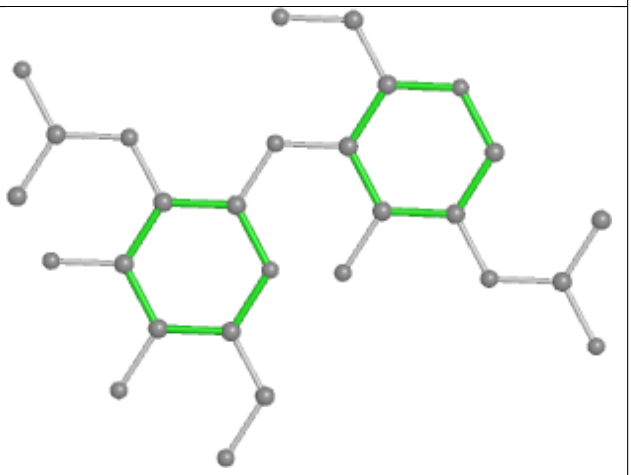
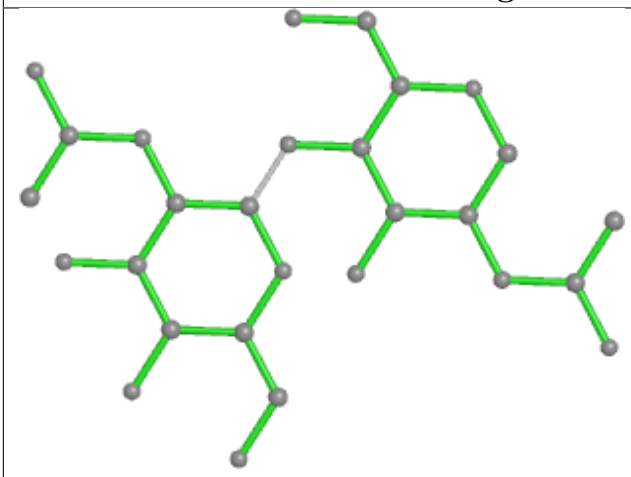
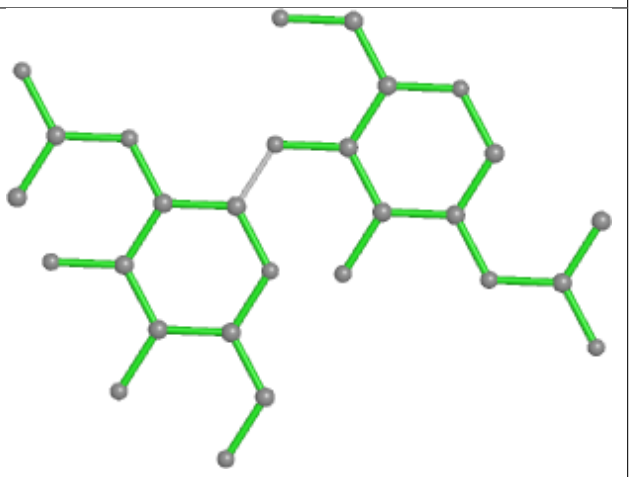


Rings

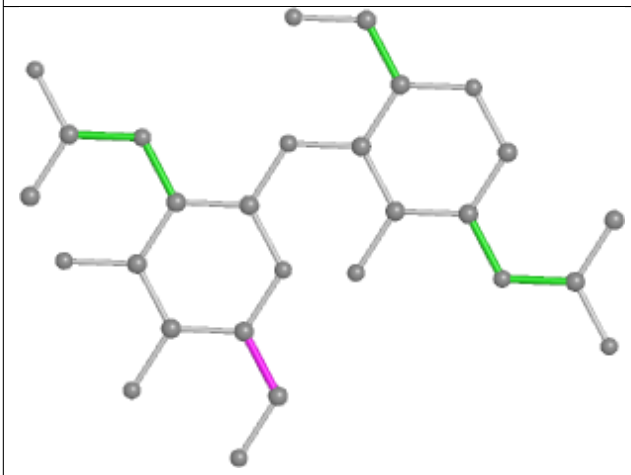
Oligosaccharide Chain L



Bond lengths



Bond angles



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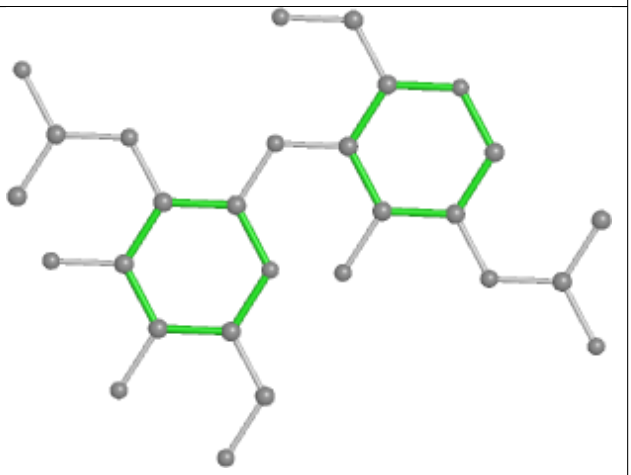
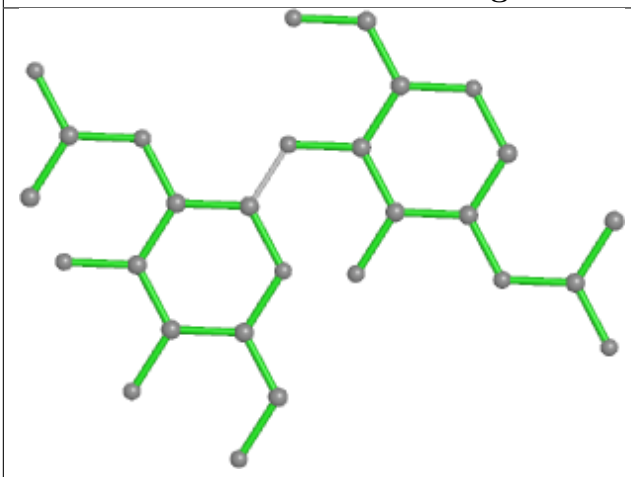


Rings

Oligosaccharide Chain M



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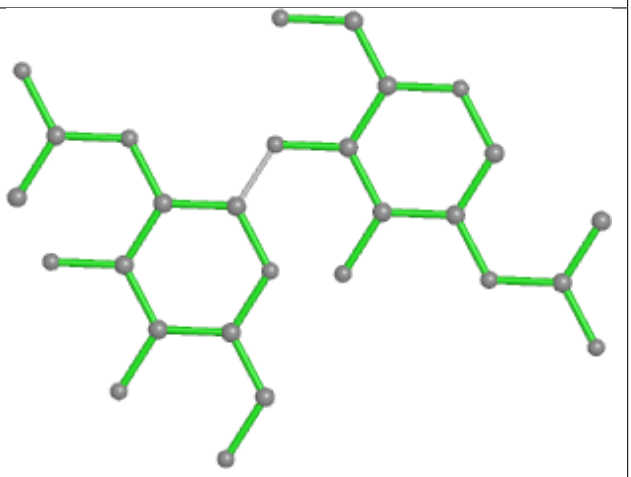
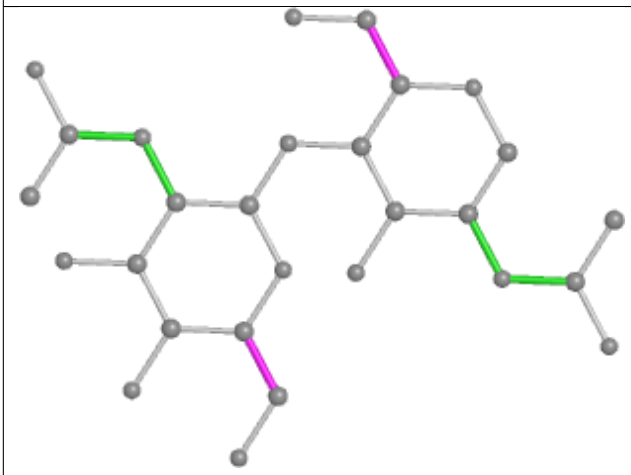


Bond angles



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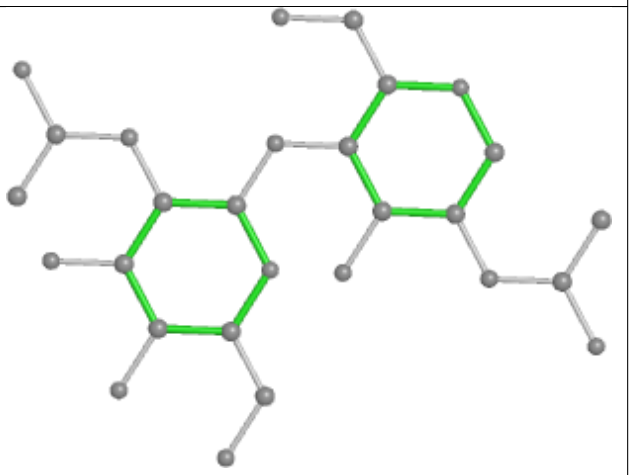
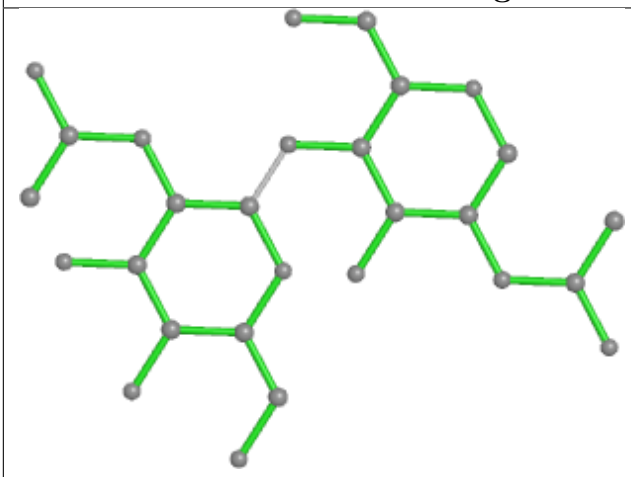


Rings

Oligosaccharide Chain N



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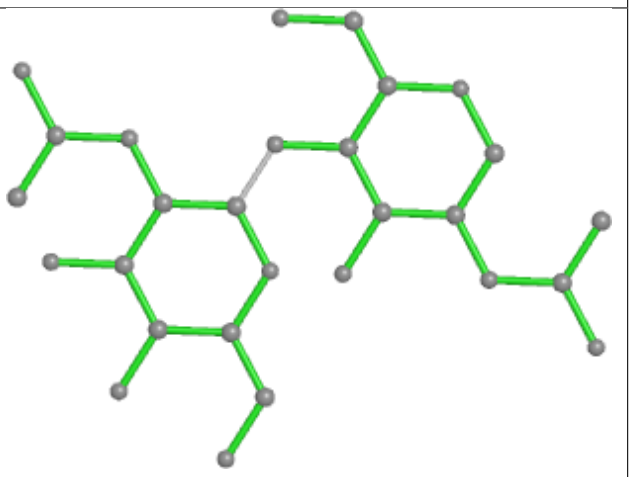
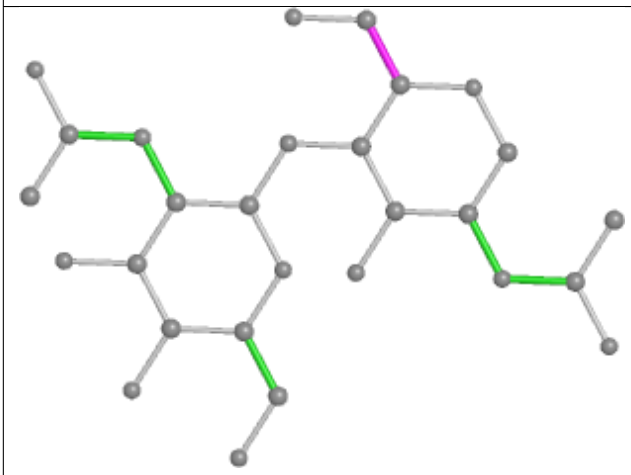


Bond angles



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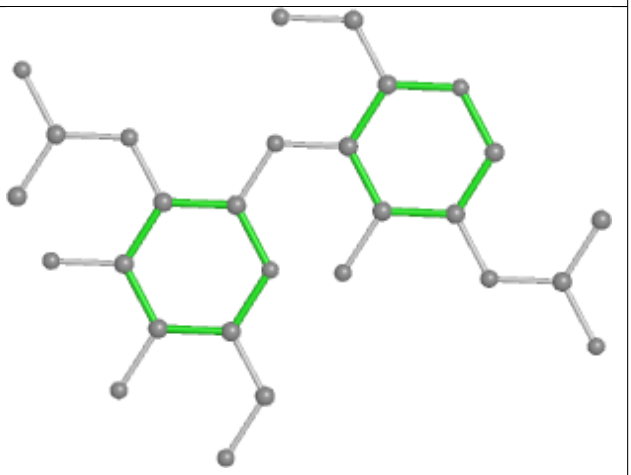
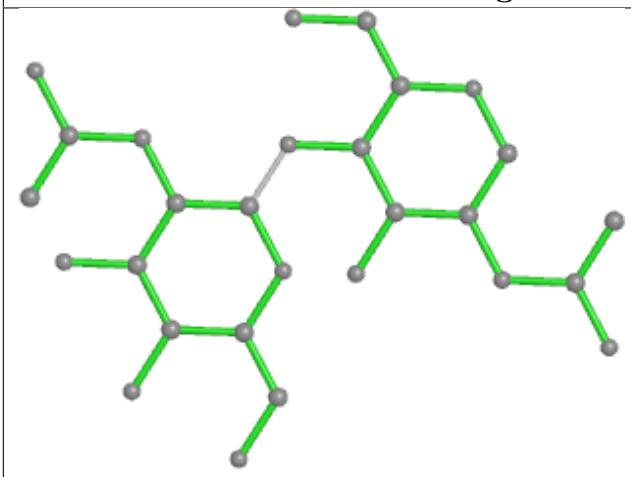


Rings

Oligosaccharide Chain P



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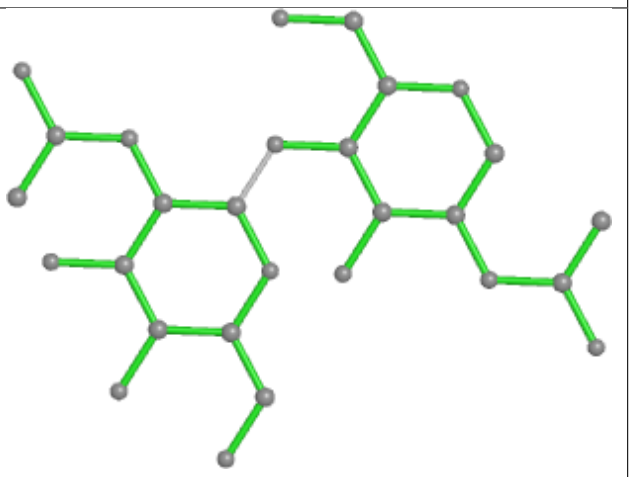

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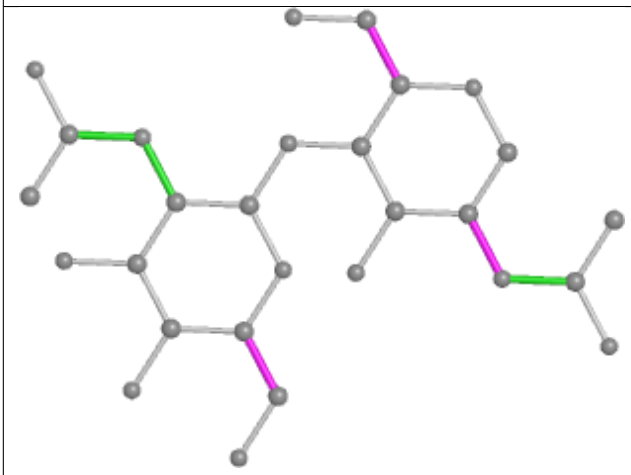

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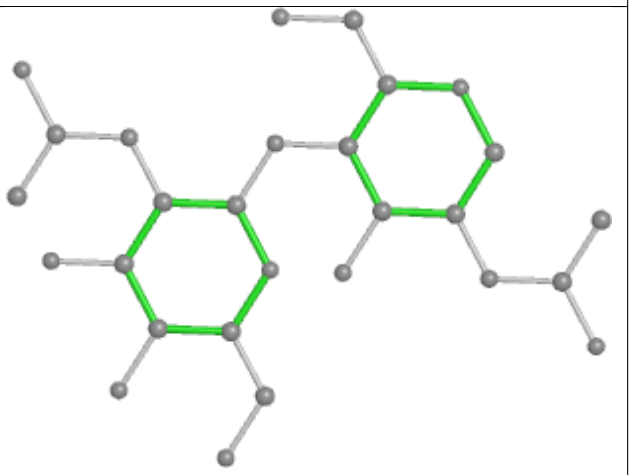
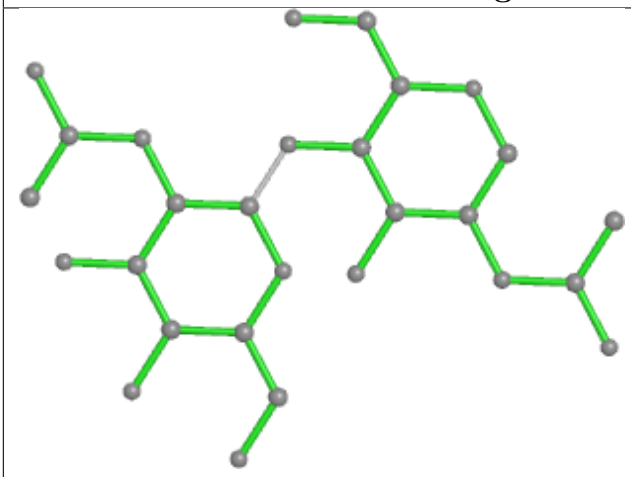


Rings

Oligosaccharide Chain Q



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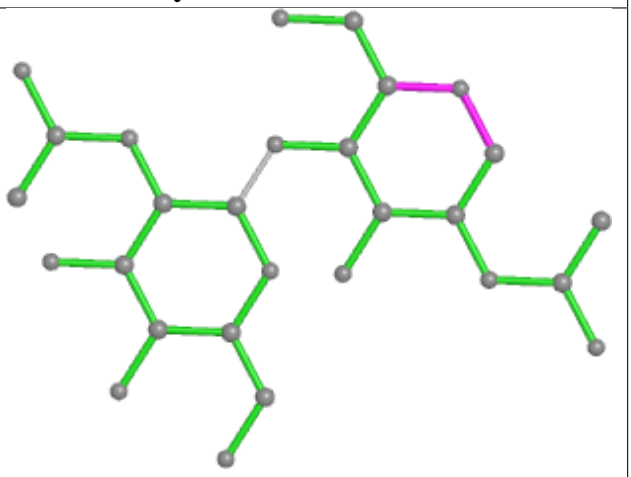

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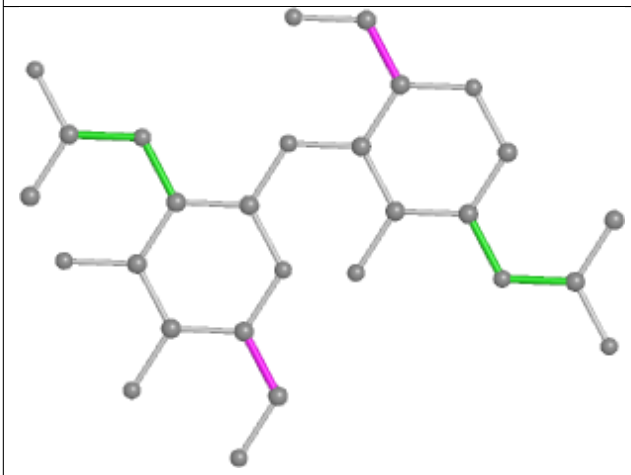

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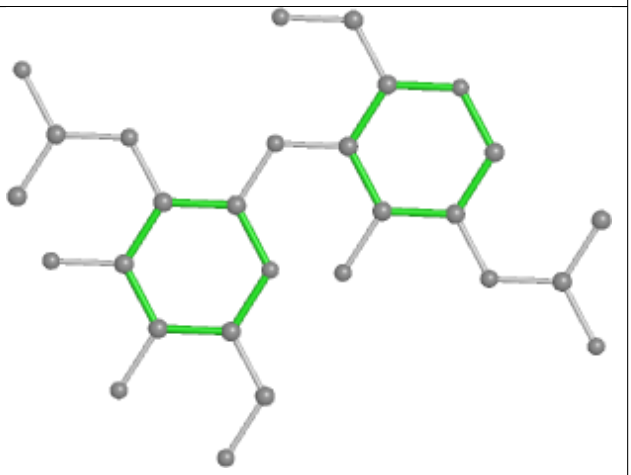
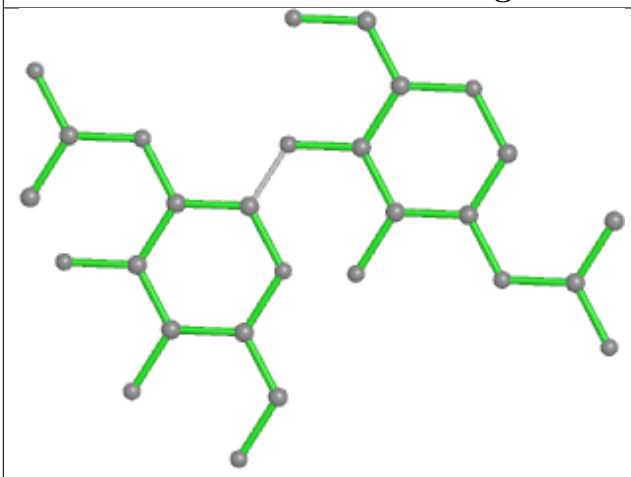


Rings

Oligosaccharide Chain R



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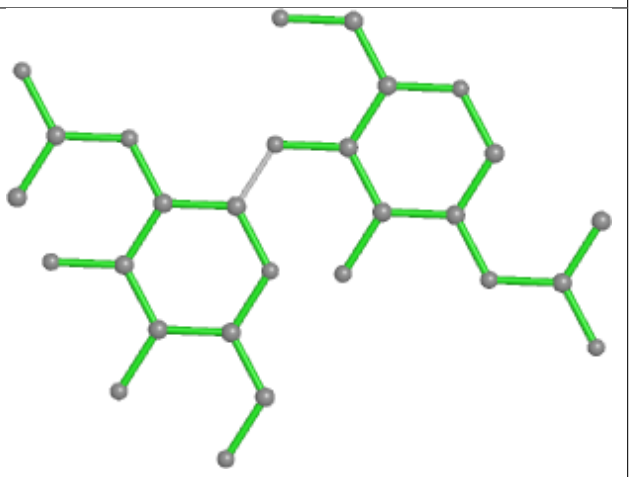

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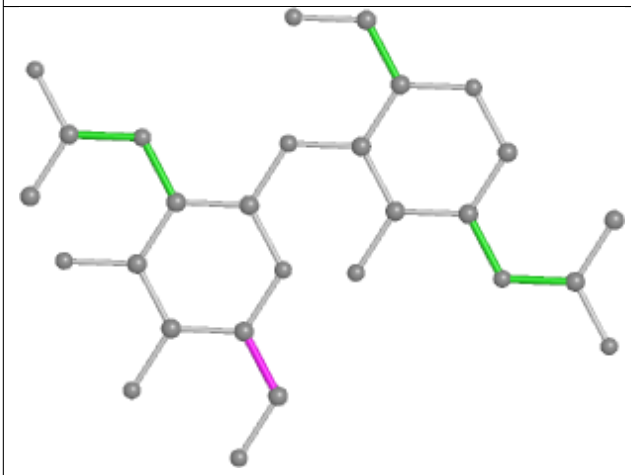

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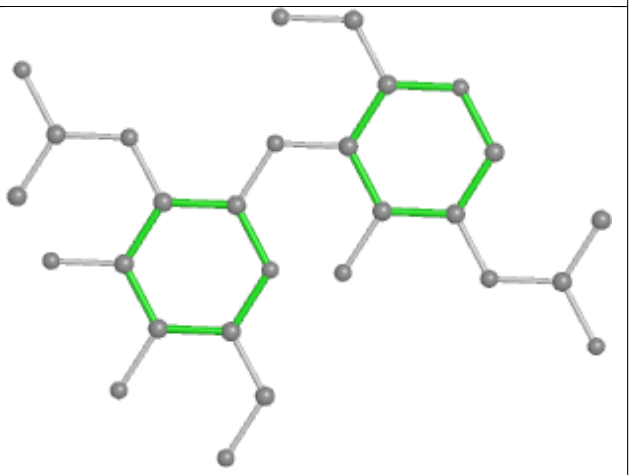
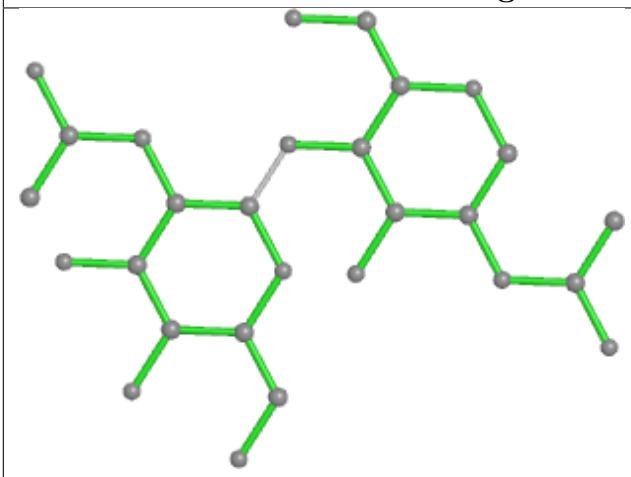


Rings

Oligosaccharide Chain S



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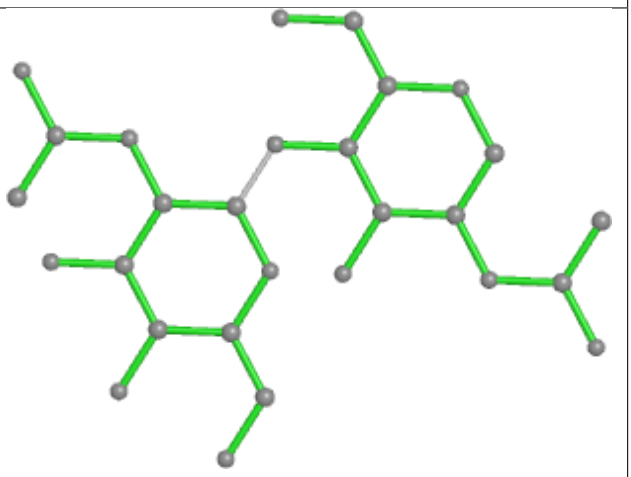

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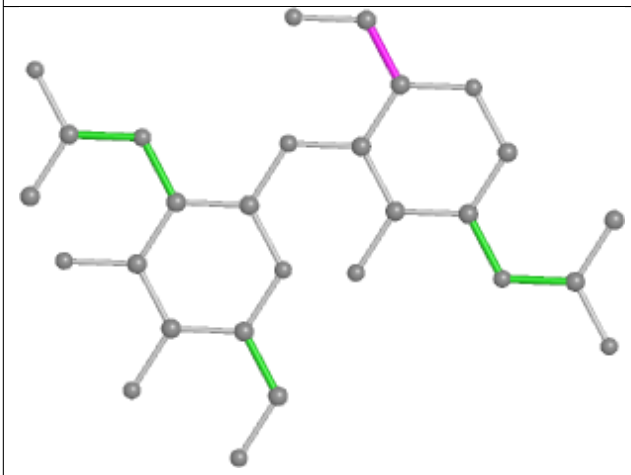

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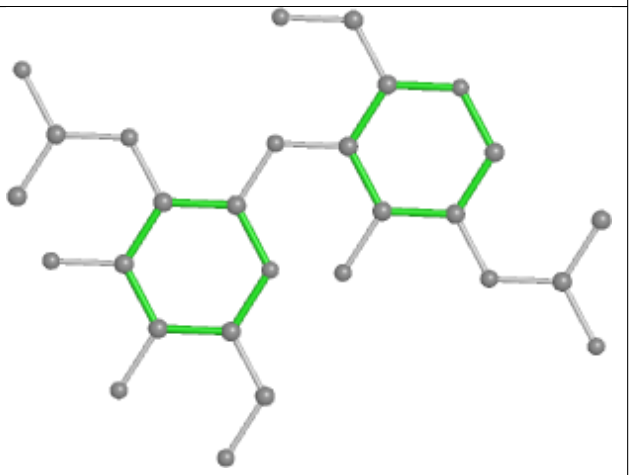
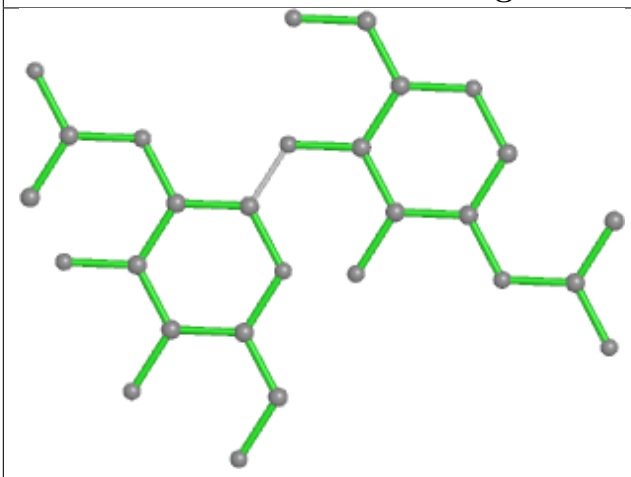


Rings

Oligosaccharide Chain T



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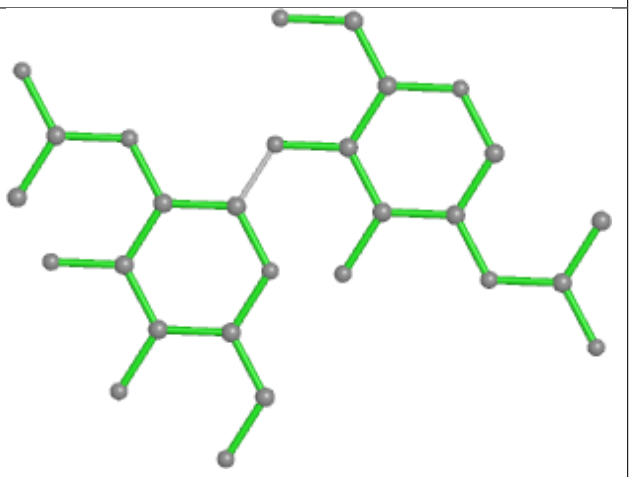

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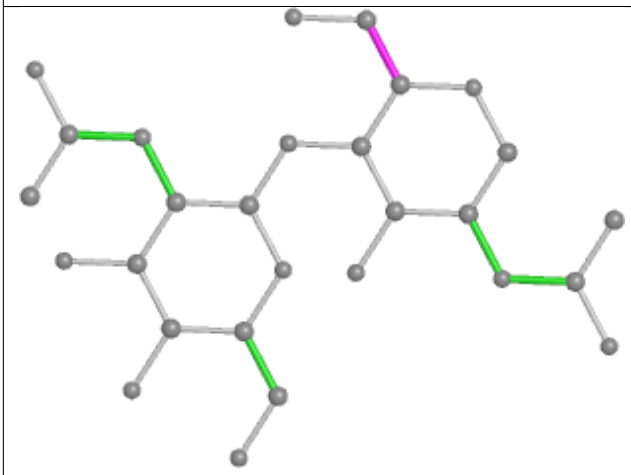

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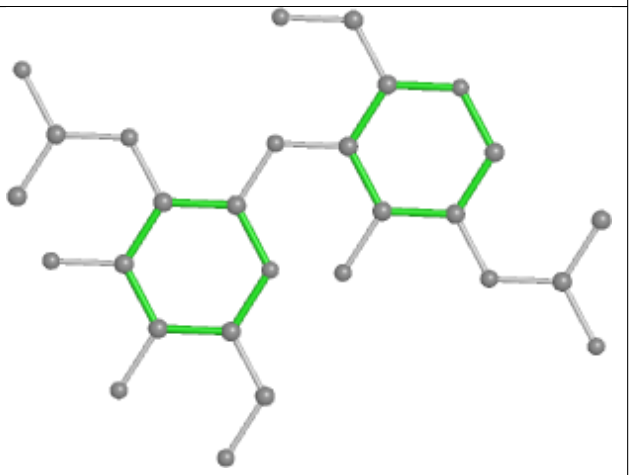
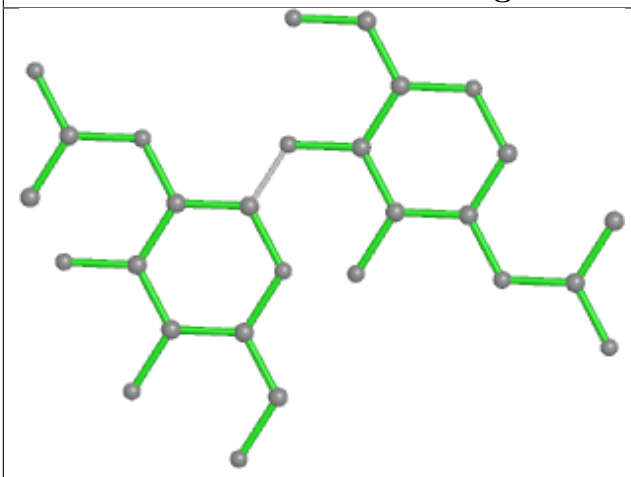


Rings

Oligosaccharide Chain U



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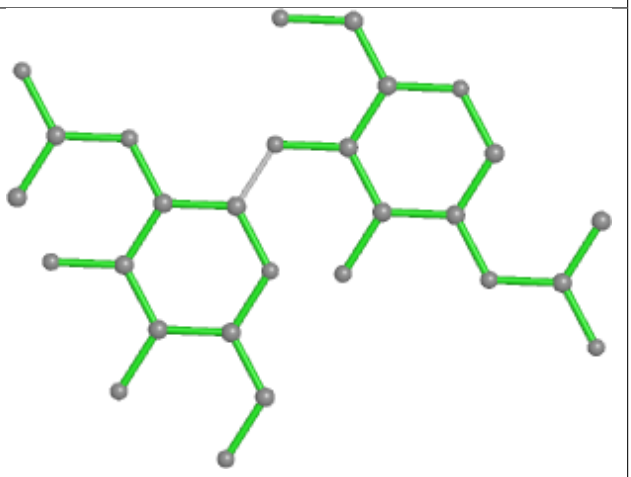

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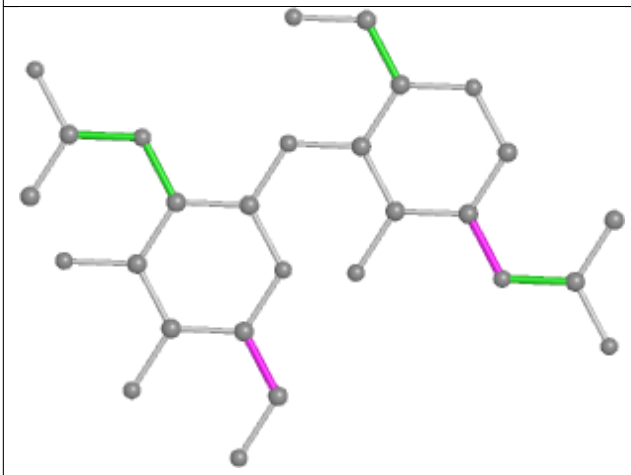

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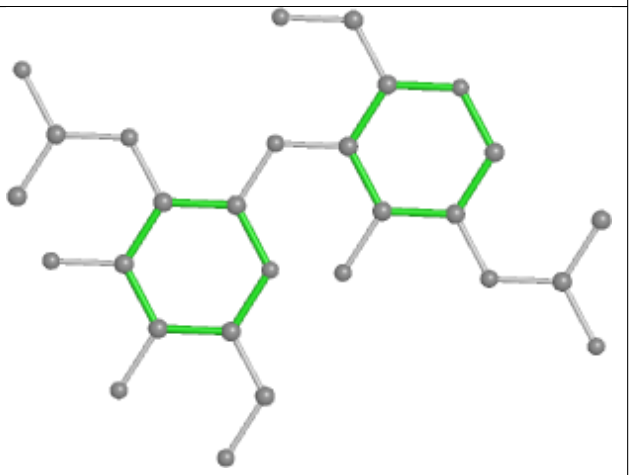
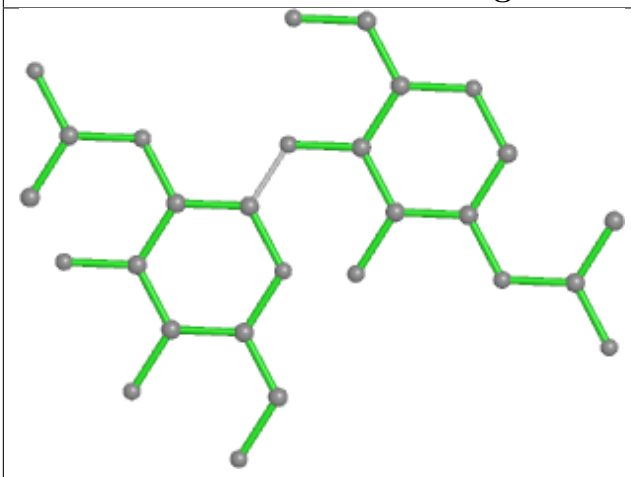


Rings

Oligosaccharide Chain V



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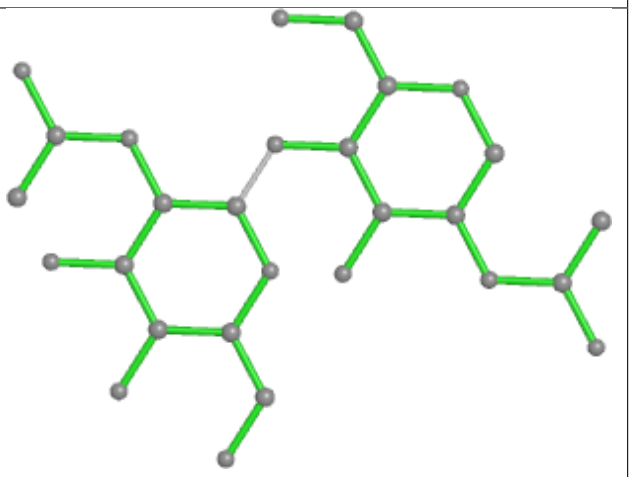
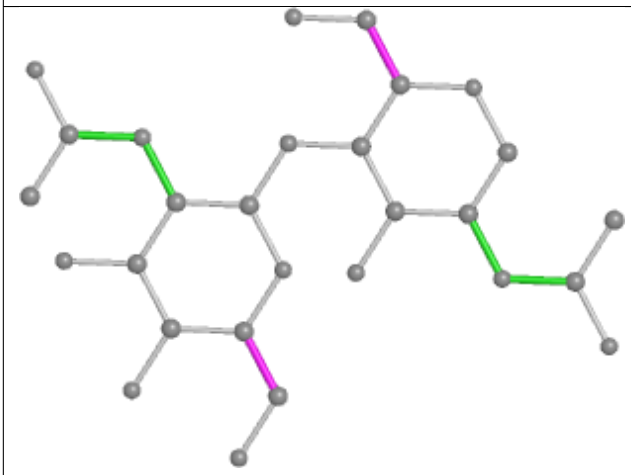


Bond angles



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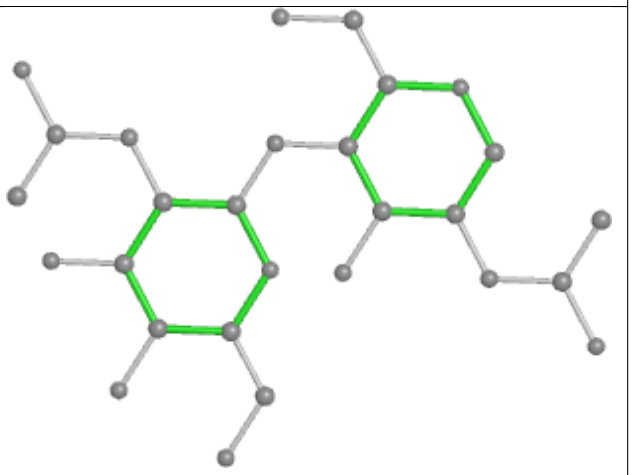
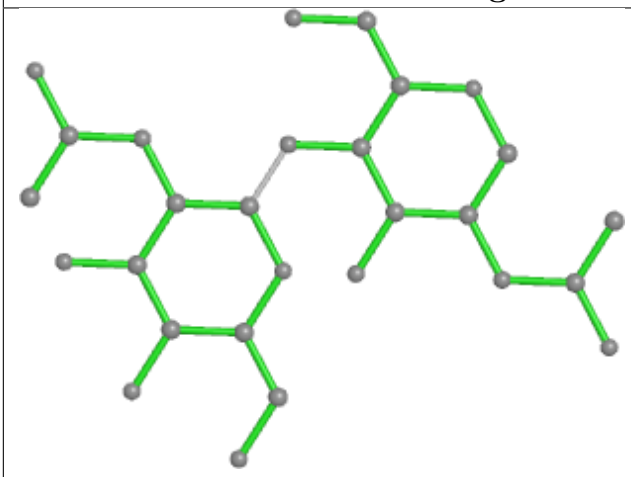


Rings

Oligosaccharide Chain W



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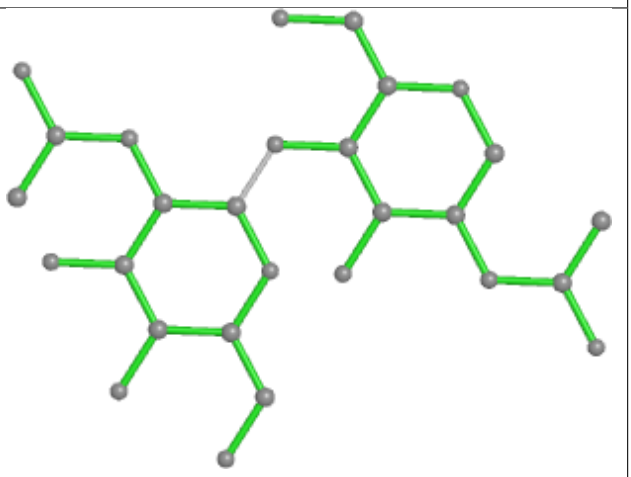
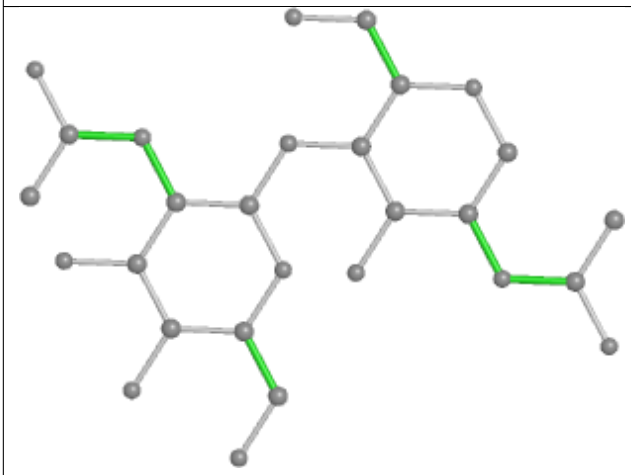


Bond angles



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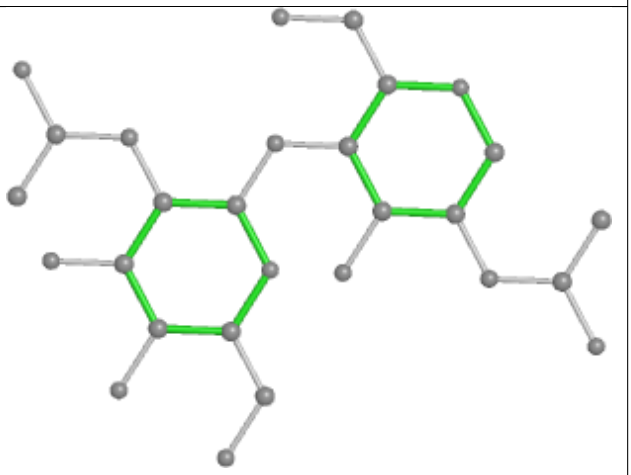
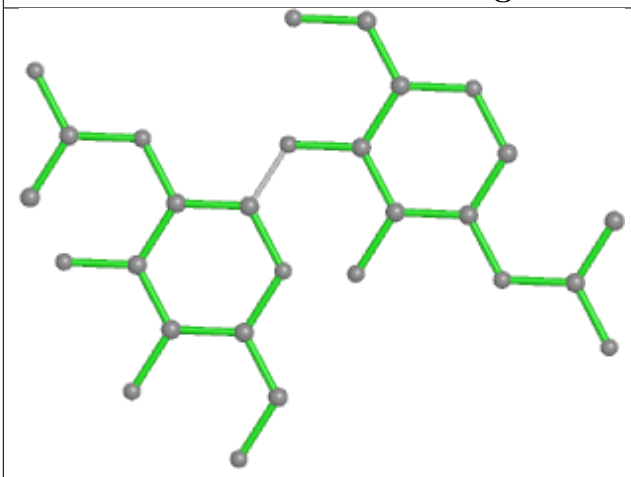


Rings

Oligosaccharide Chain X



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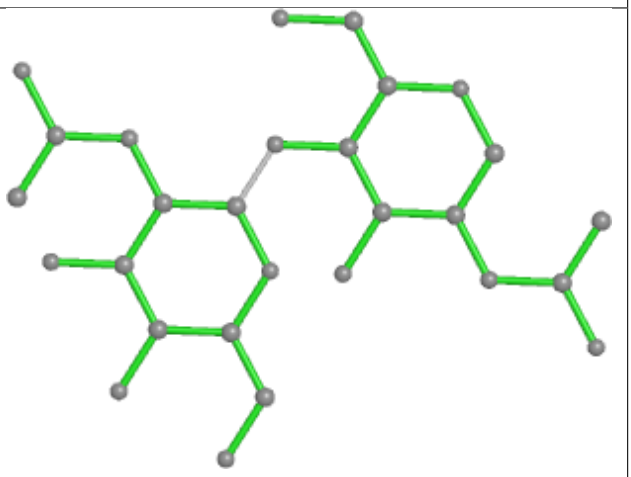
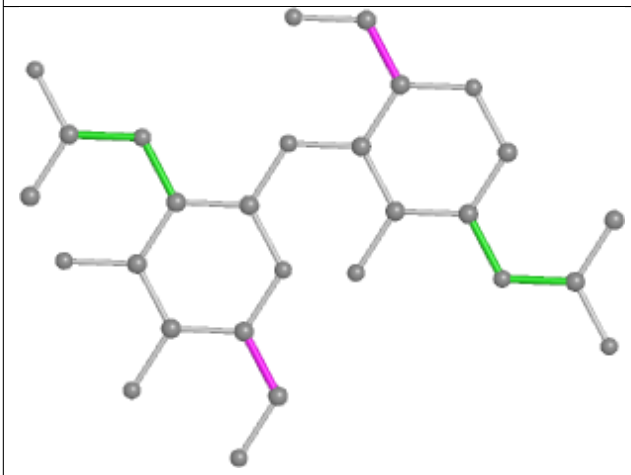
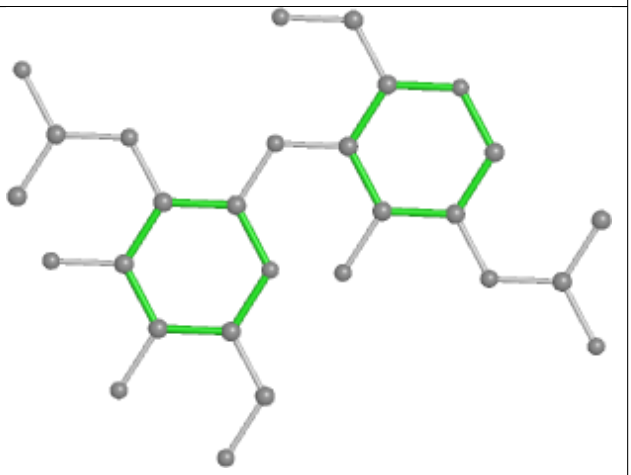


Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

14 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	NAG	B	1304	1	14,14,15	0.26	0	17,19,21	0.49	0
5	NAG	C	1304	1	14,14,15	0.23	0	17,19,21	0.43	0
5	NAG	A	1302	1	14,14,15	0.24	0	17,19,21	0.43	0
5	NAG	C	1302	1	14,14,15	0.20	0	17,19,21	0.46	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	NAG	A	1305	1	14,14,15	0.21	0	17,19,21	0.43	0
5	NAG	C	1303	1	14,14,15	0.30	0	17,19,21	0.45	0
5	NAG	A	1301	1	14,14,15	0.18	0	17,19,21	0.45	0
5	NAG	C	1301	1	14,14,15	0.20	0	17,19,21	0.37	0
5	NAG	A	1304	1	14,14,15	0.25	0	17,19,21	0.52	0
5	NAG	B	1301	1	14,14,15	0.22	0	17,19,21	0.41	0
5	NAG	B	1303	1	14,14,15	0.29	0	17,19,21	0.39	0
5	NAG	B	1302	1	14,14,15	0.19	0	17,19,21	0.38	0
5	NAG	A	1303	1	14,14,15	0.33	0	17,19,21	0.35	0
5	NAG	C	1305	1	14,14,15	0.20	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
5	NAG	B	1304	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1304	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1302	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1302	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1305	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1303	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1301	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1301	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1304	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1301	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1303	1	-	1/6/23/26	0/1/1/1
5	NAG	B	1302	1	-	2/6/23/26	0/1/1/1
5	NAG	A	1303	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1305	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1303	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	C	1302	NAG	O5-C5-C6-O6
5	B	1302	NAG	O5-C5-C6-O6
5	A	1303	NAG	C4-C5-C6-O6
5	C	1301	NAG	O5-C5-C6-O6
5	C	1301	NAG	C4-C5-C6-O6
5	C	1302	NAG	C4-C5-C6-O6
5	B	1304	NAG	O5-C5-C6-O6
5	C	1304	NAG	O5-C5-C6-O6
5	A	1304	NAG	C4-C5-C6-O6
5	C	1304	NAG	C4-C5-C6-O6
5	B	1304	NAG	C4-C5-C6-O6
5	B	1302	NAG	C4-C5-C6-O6
5	A	1304	NAG	O5-C5-C6-O6
5	C	1303	NAG	C1-C2-N2-C7
5	B	1303	NAG	C1-C2-N2-C7

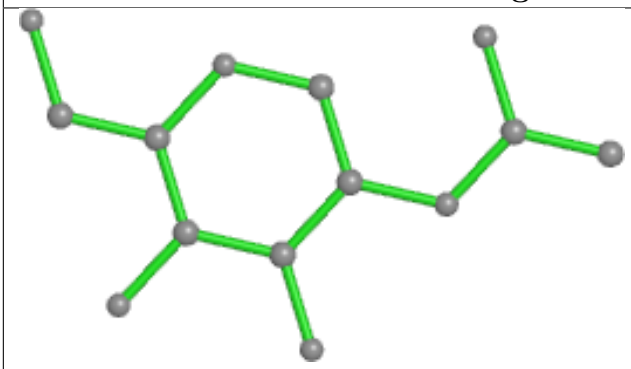
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1303	NAG	1	0
5	B	1302	NAG	1	0
5	A	1303	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 then 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 1304



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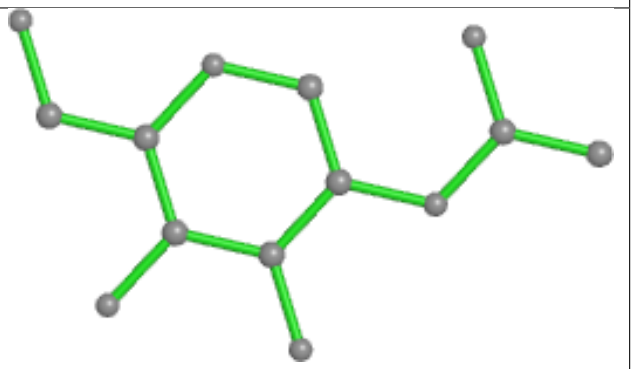
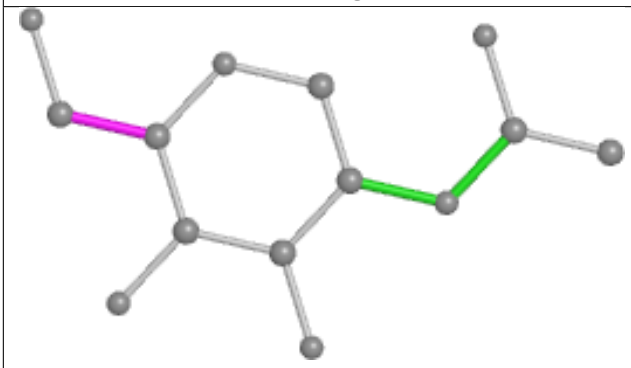


Bond angles



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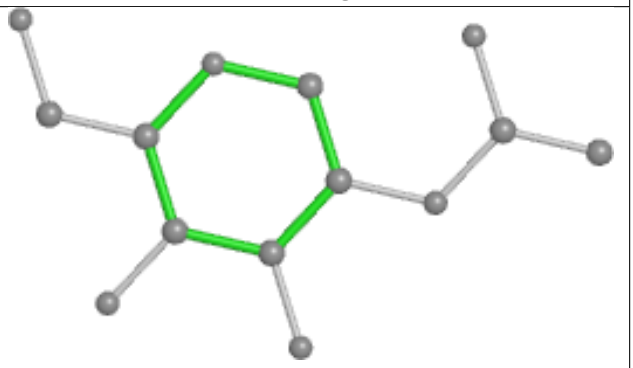
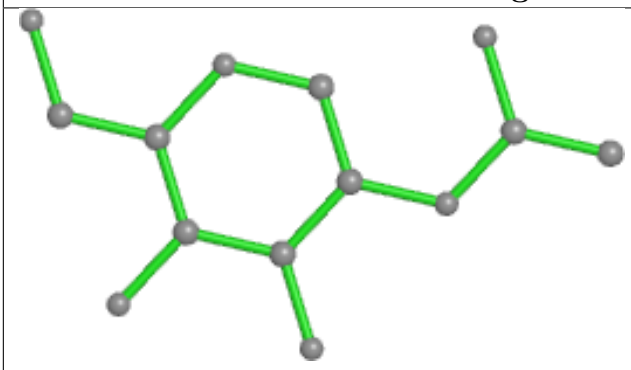


Rings

Ligand NAG C 1304



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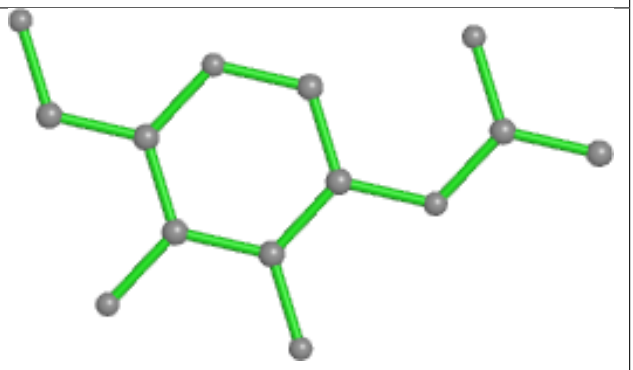

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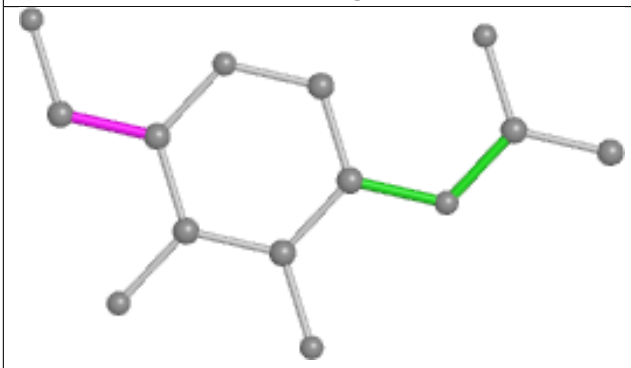

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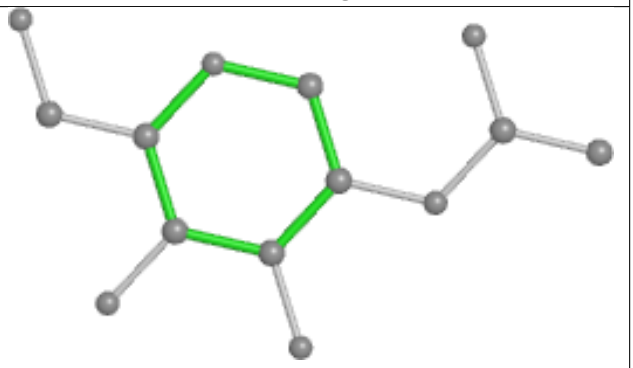
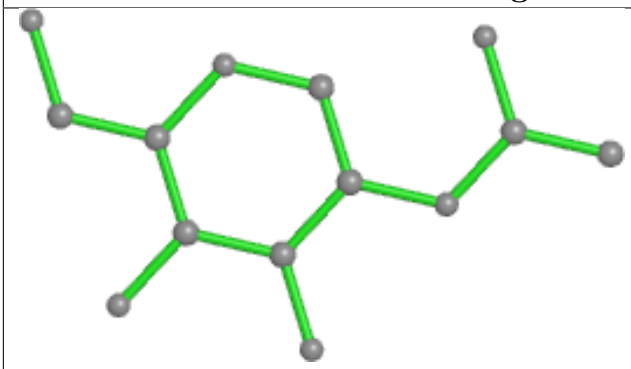


Rings

Ligand NAG A 1302



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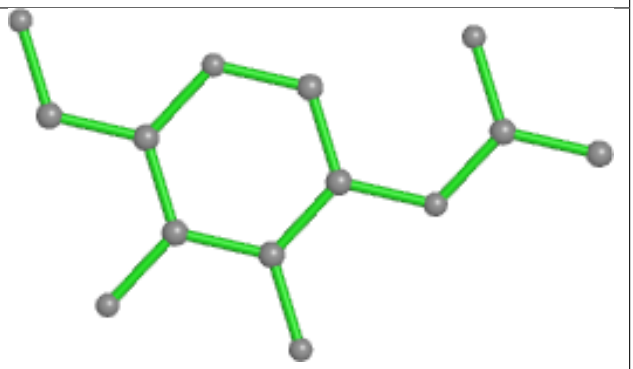

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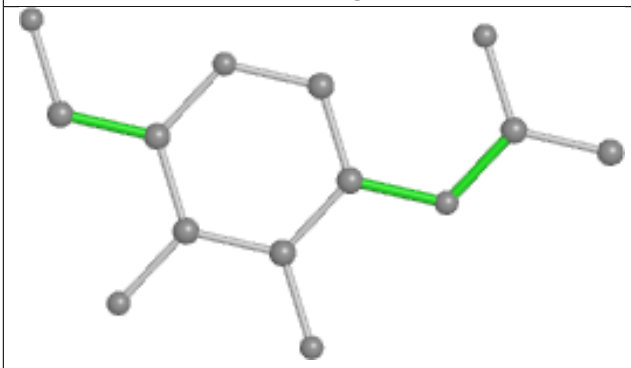

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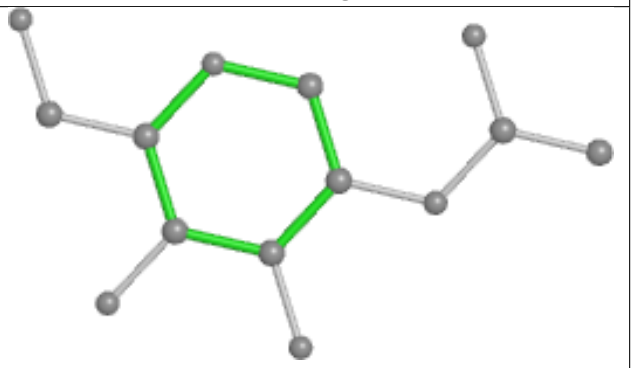

Rings

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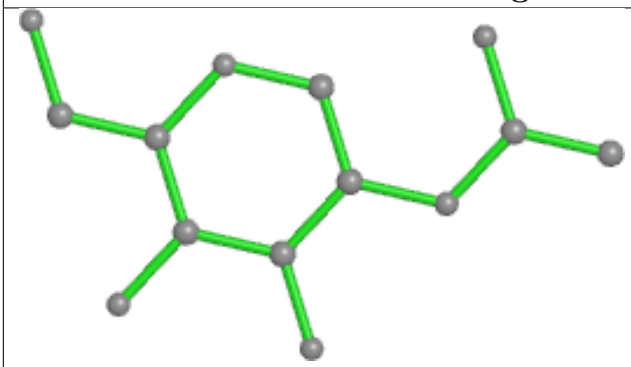

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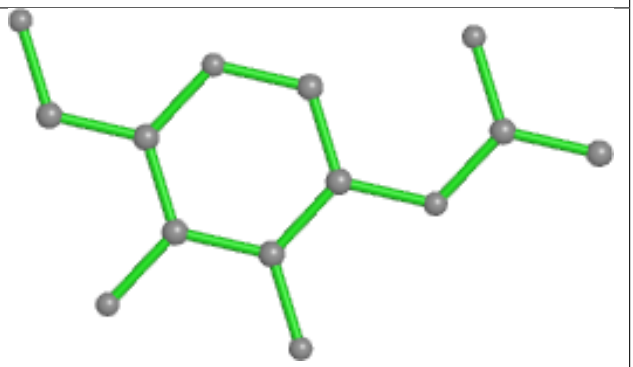

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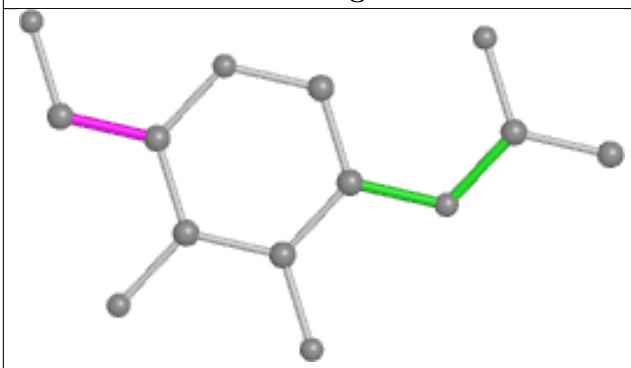

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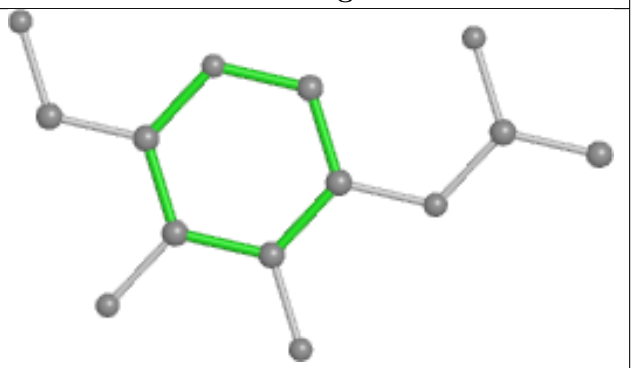
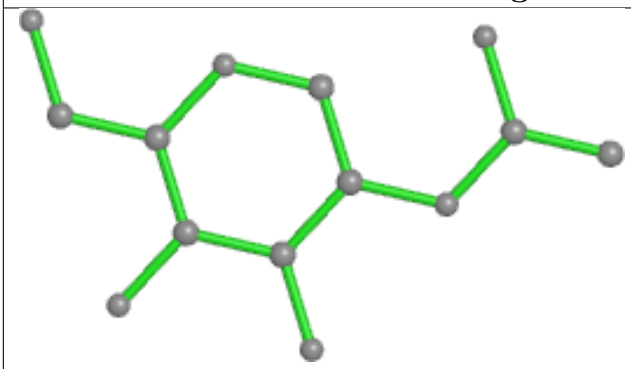


Rings

Ligand NAG A 1305



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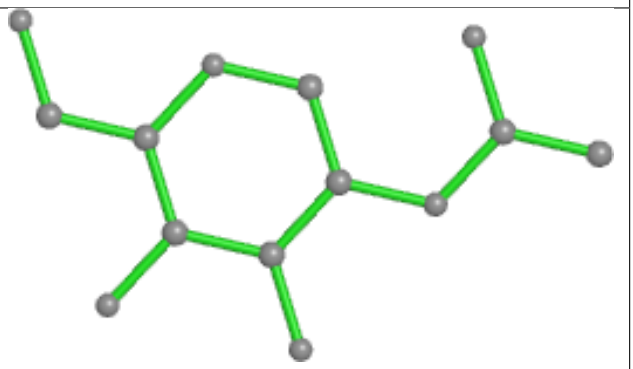

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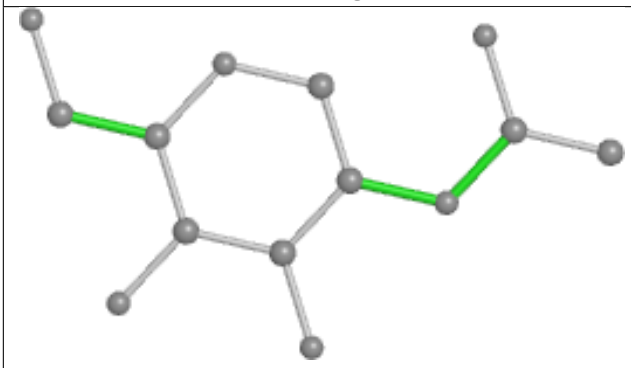

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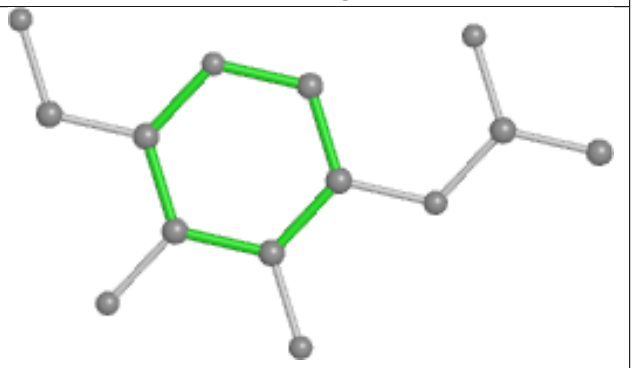

Rings

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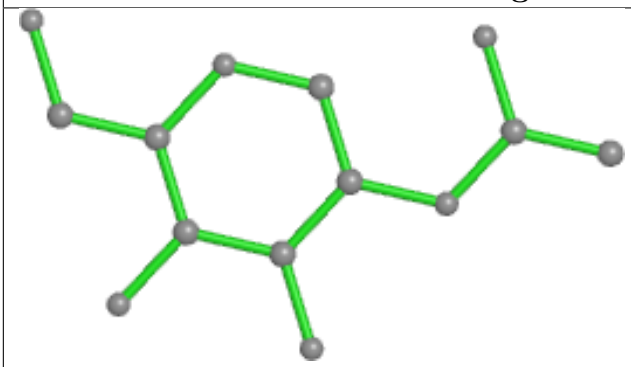

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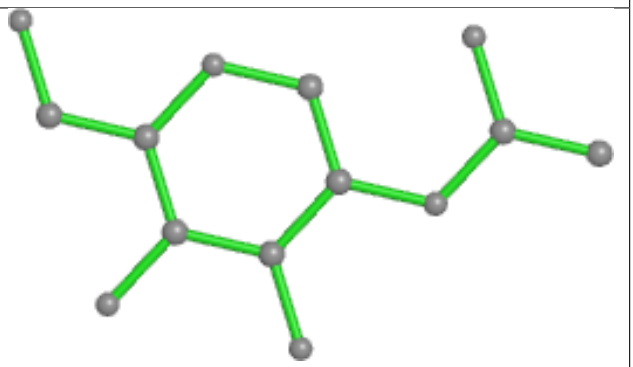

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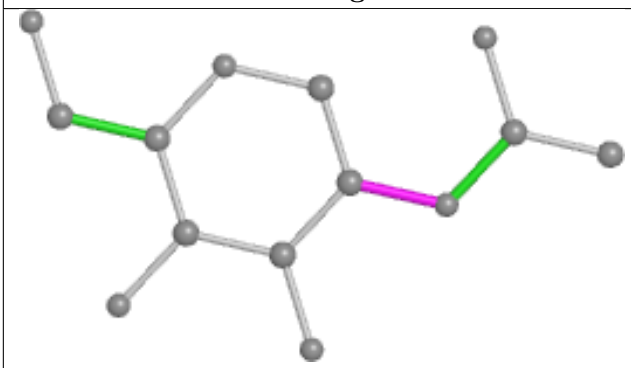

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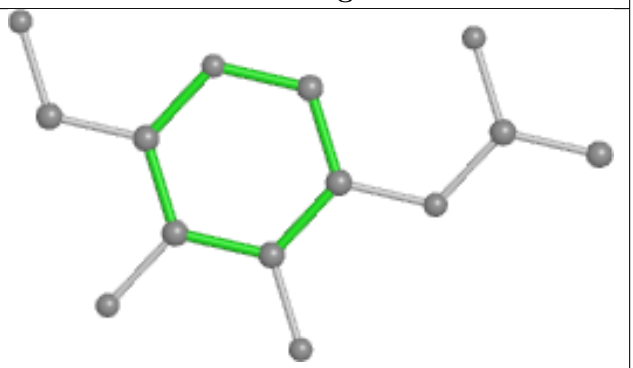
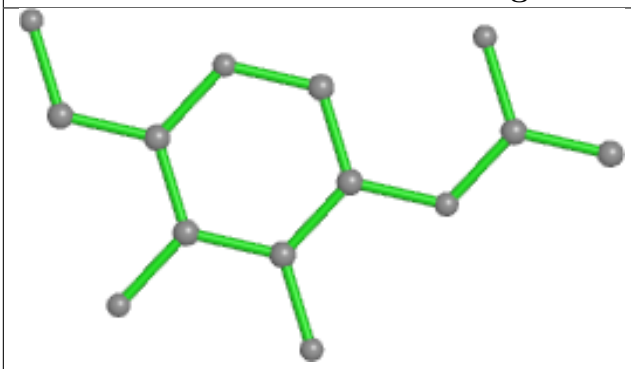


Rings

Ligand NAG A 1301



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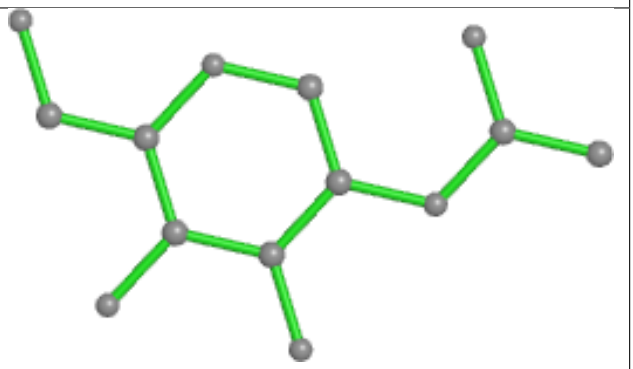

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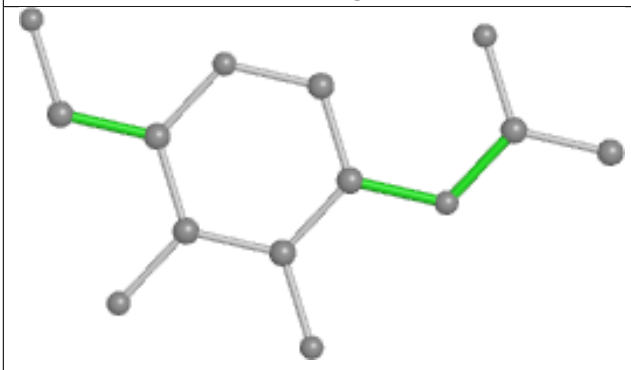

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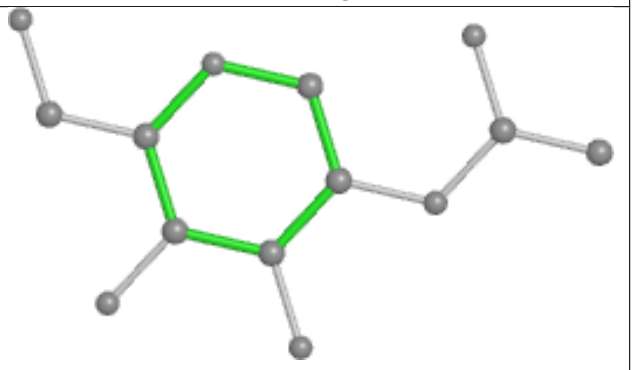

Rings

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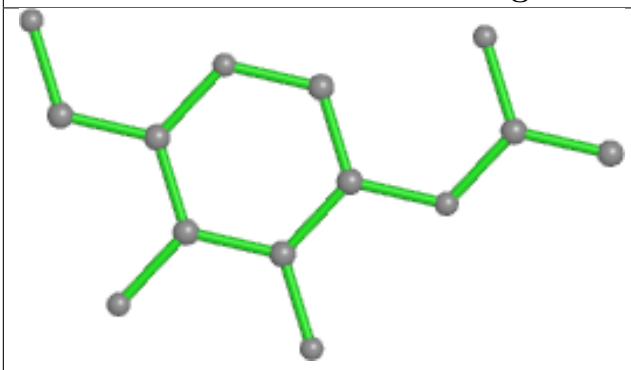

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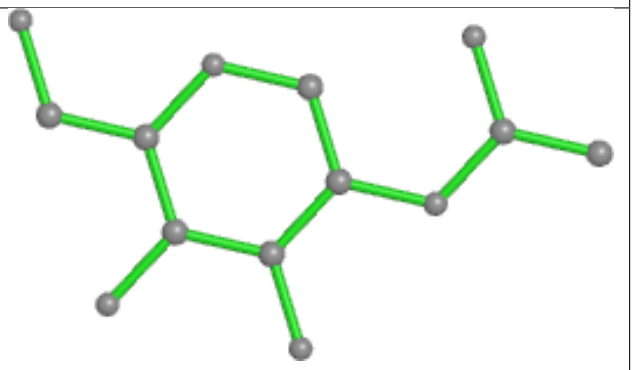
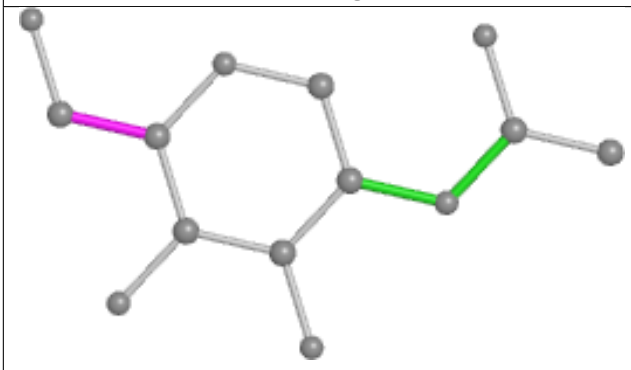
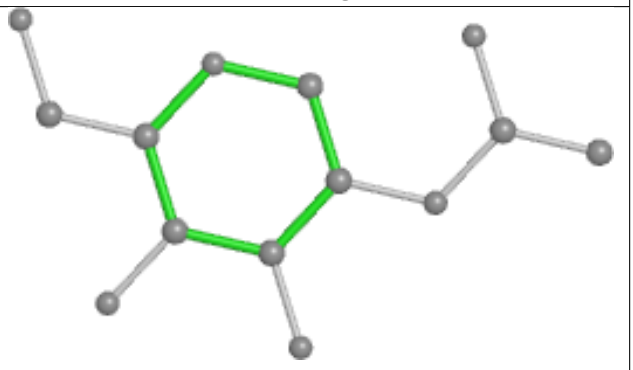


Bond angles

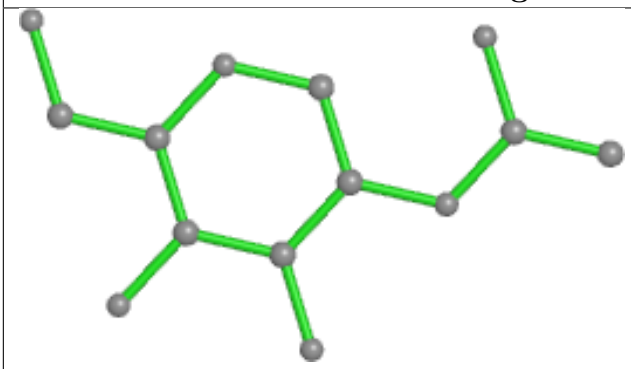


Torsions

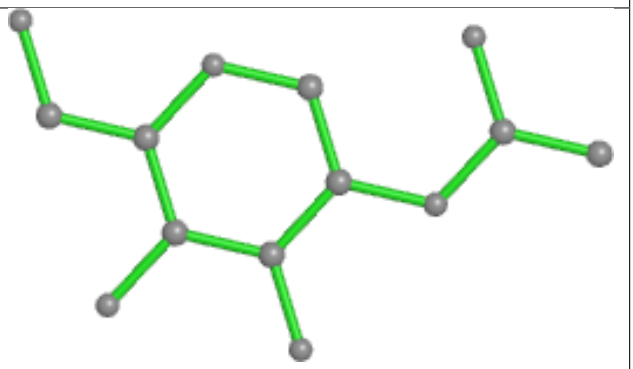


Rings

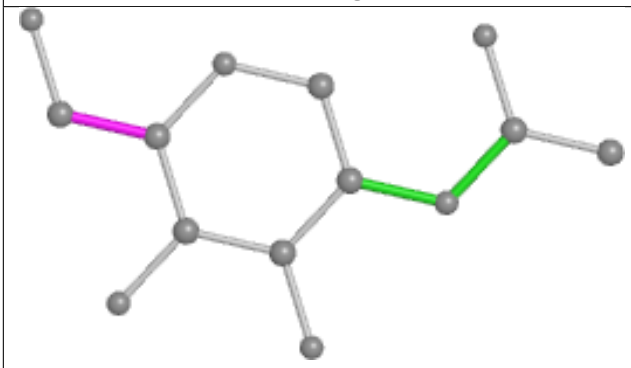
Ligand NAG A 1304



Bond lengths



Bond angles



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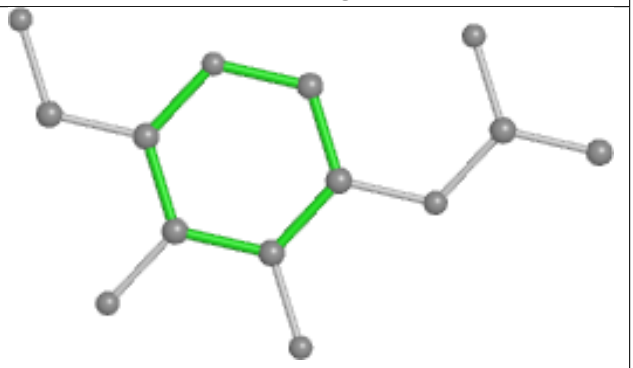

Rings

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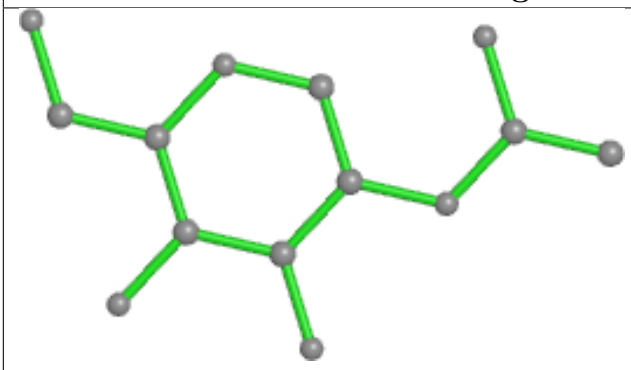

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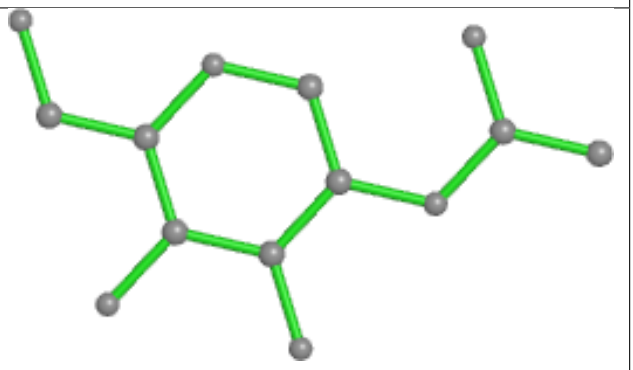
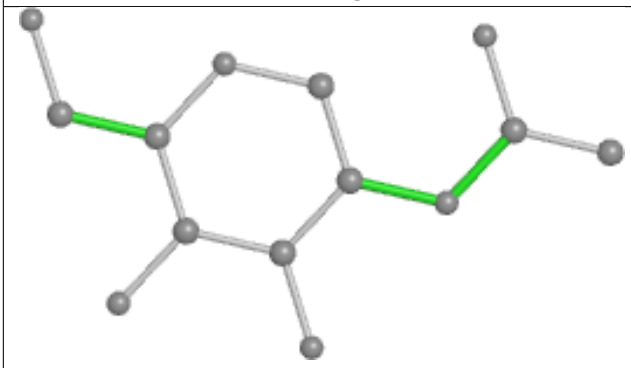
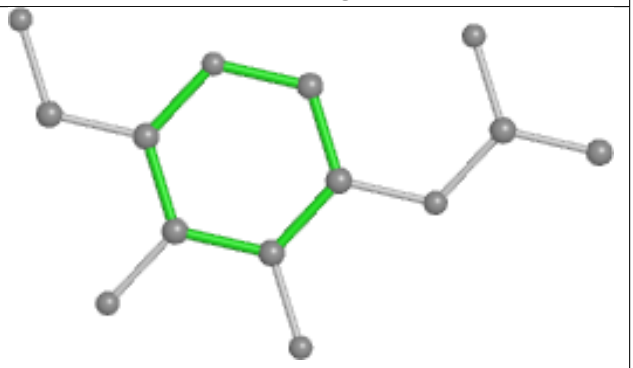


Bond angles

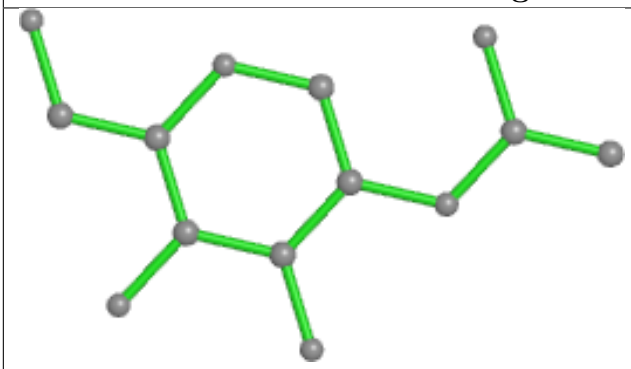


Torsions

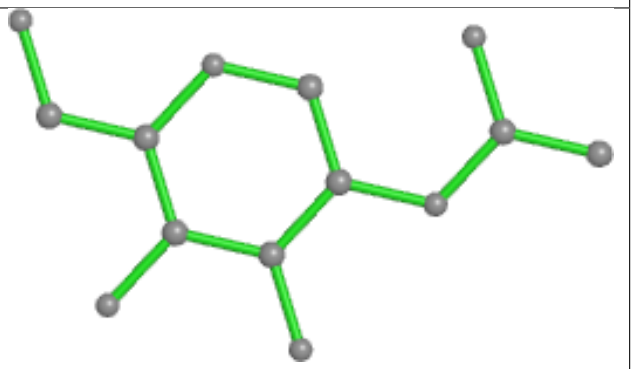


Rings

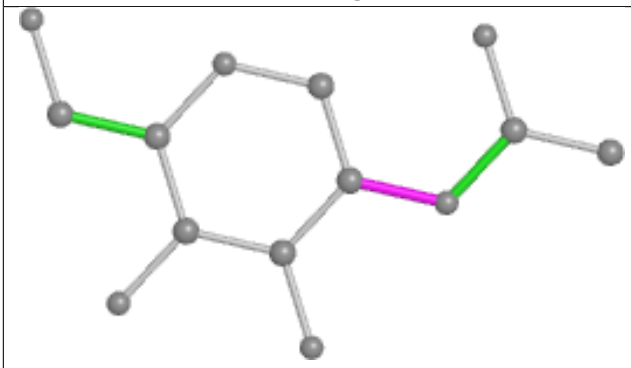
Ligand NAG B 1303



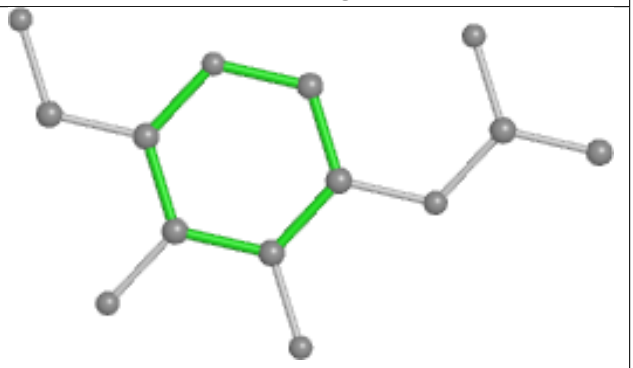
Bond lengths



Bond angles

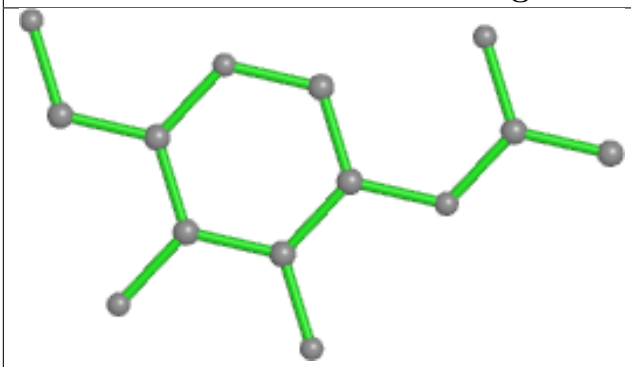


Torsions

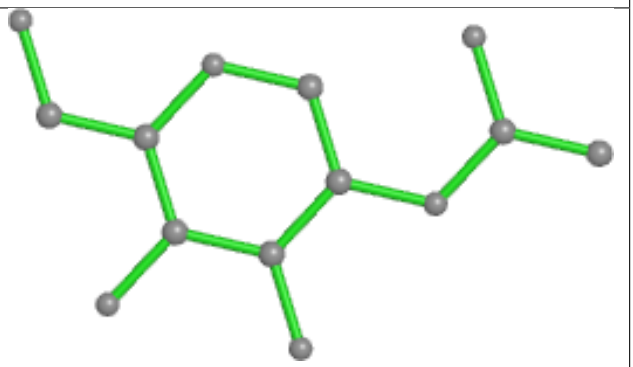


Rings

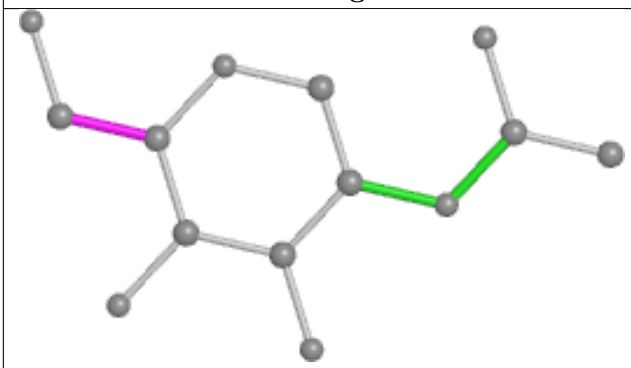
Ligand NAG B 1302



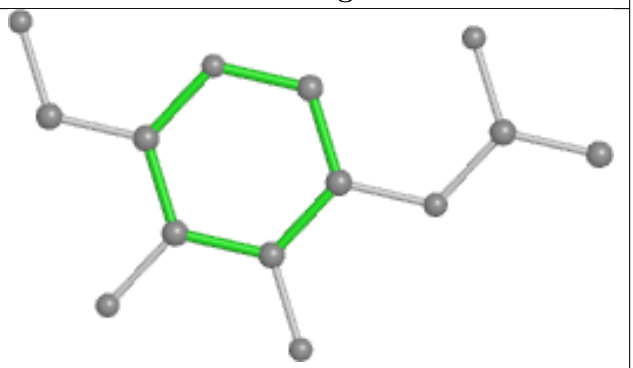
Bond lengths



Bond angles

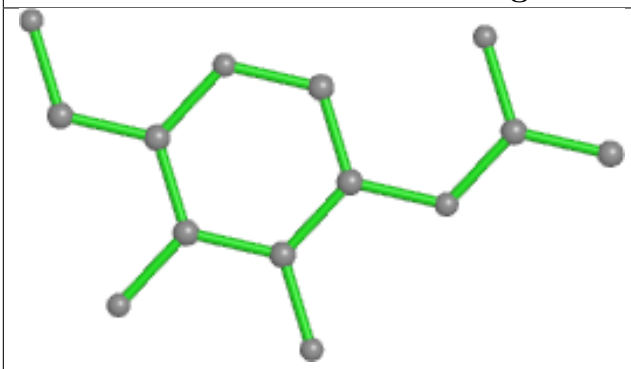


Torsions

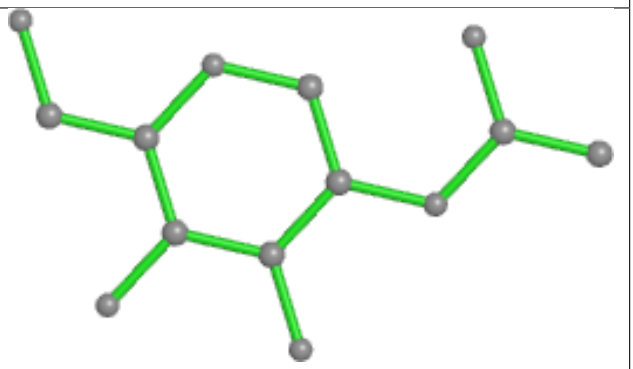


Rings

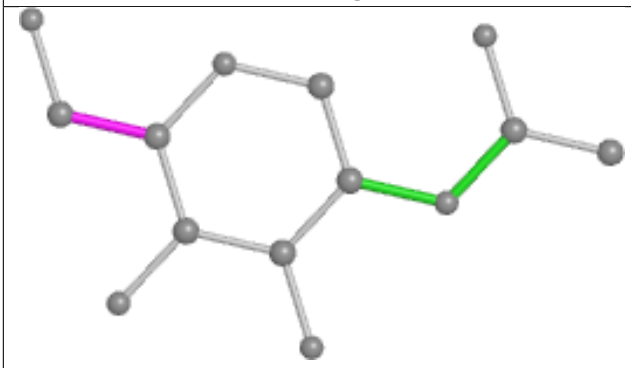
Ligand NAG A 1303



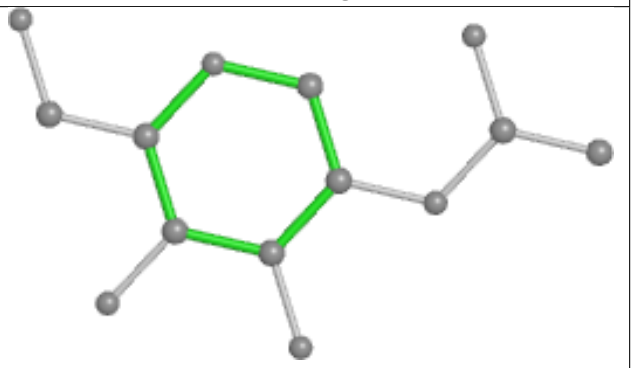
Bond lengths



Bond angles

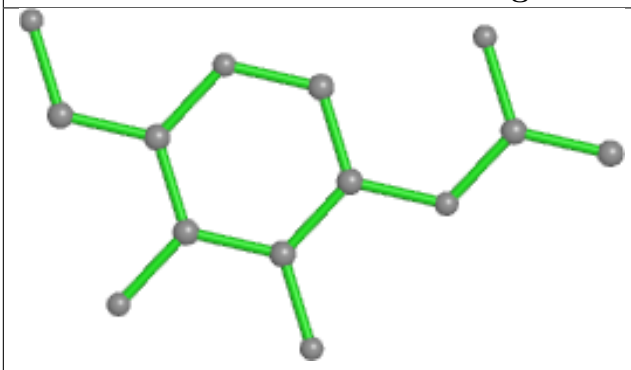


Torsions

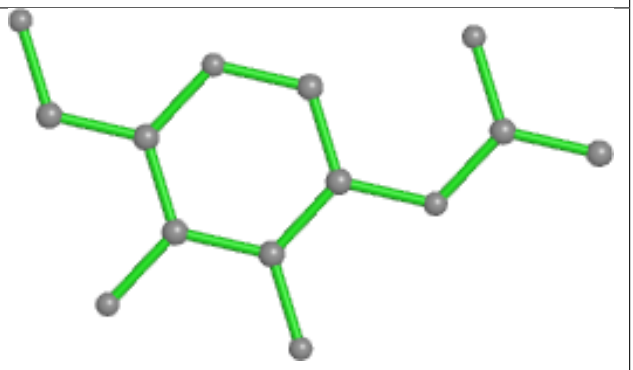


Rings

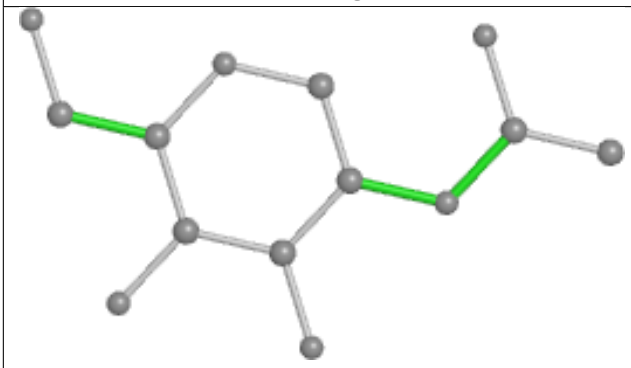
Ligand NAG C 1305



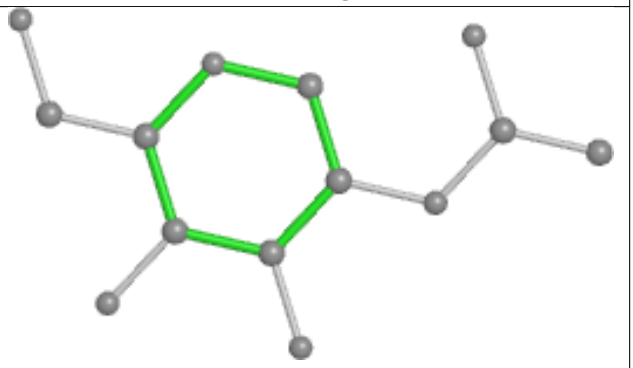
Bond lengths



Bond angles



Torsions



Rings

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.

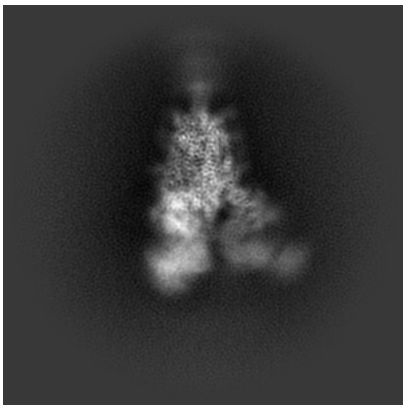
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-34408. These allow visual inspection of the internal detail of the map and identification of artifacts.

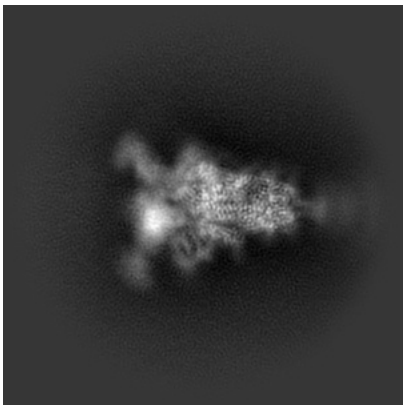
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

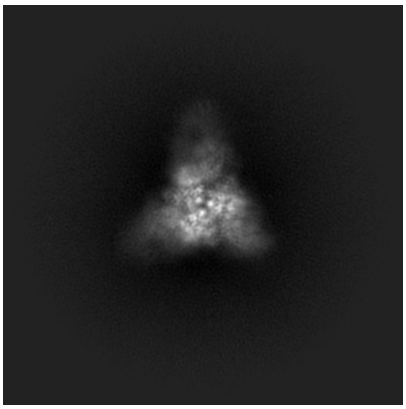
6.1.1 Primary map



X

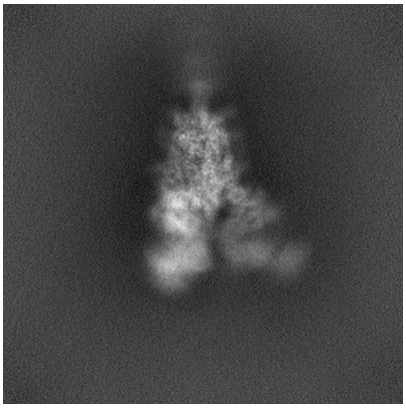


Y

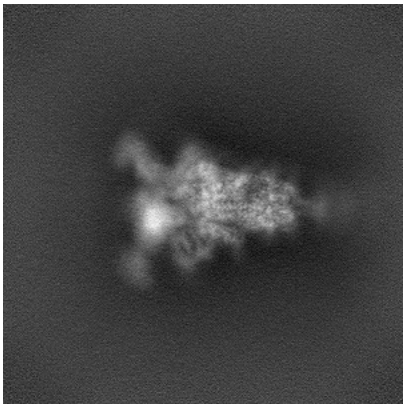


Z

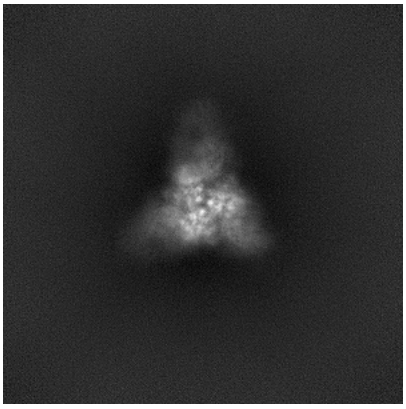
6.1.2 Raw map



X



Y

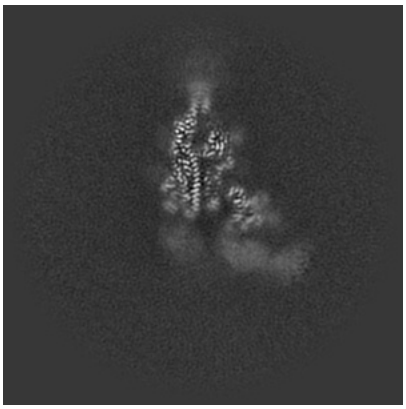


Z

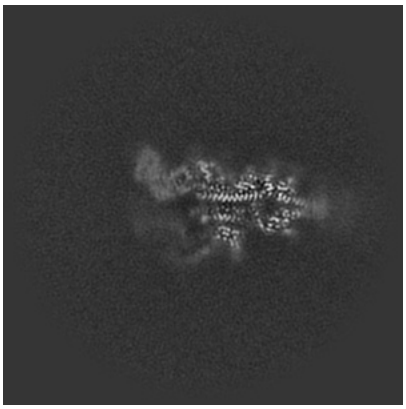
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

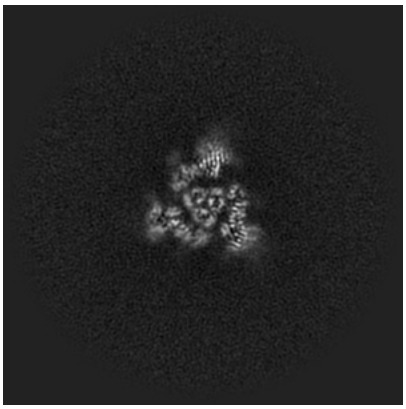
6.2.1 Primary map



X Index: 256

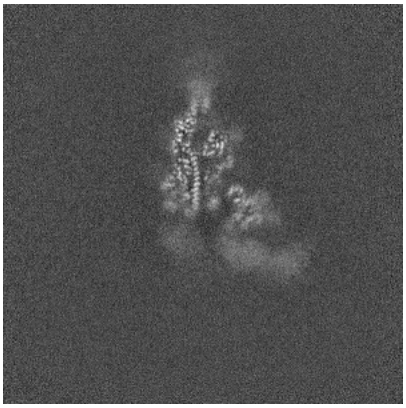


Y Index: 256

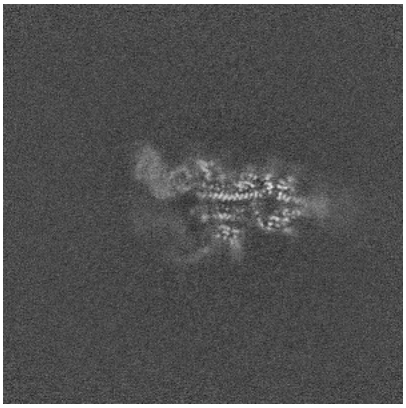


Z Index: 256

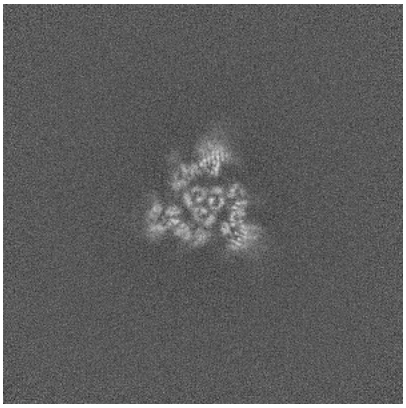
6.2.2 Raw map



X Index: 256



Y Index: 256



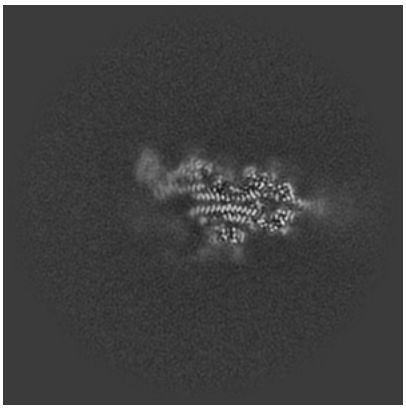
Z Index: 256

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

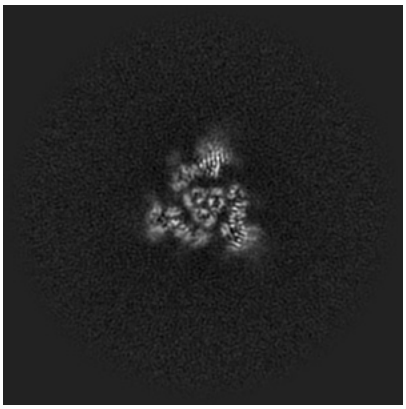
6.3 Largest variance slices (i)

6.3.1 Primary map

X Index: 251

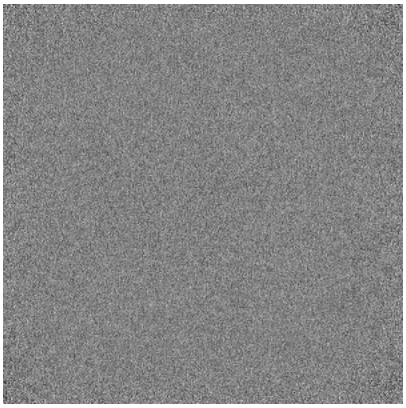


Y Index: 260

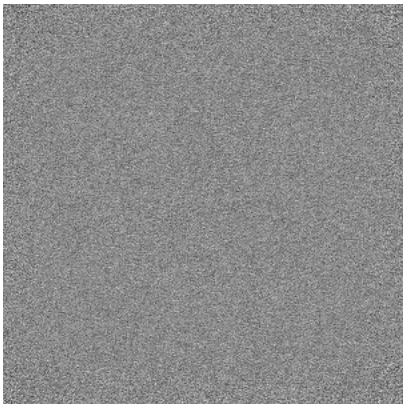


Z Index: 256

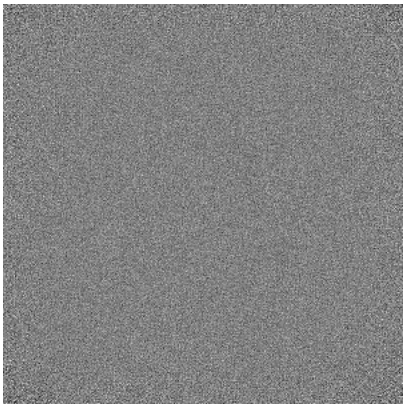
6.3.2 Raw map



X Index: 0



Y Index: 0

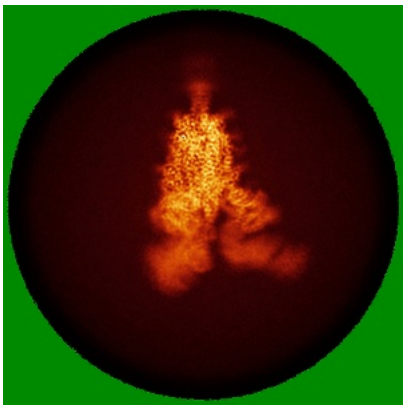


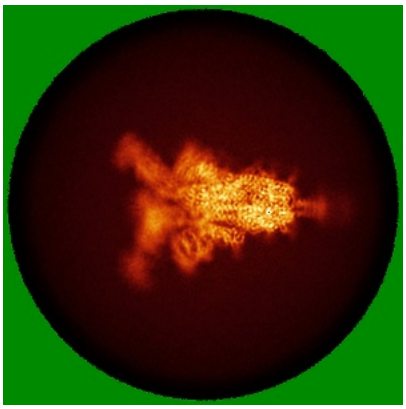
Z Index: 0

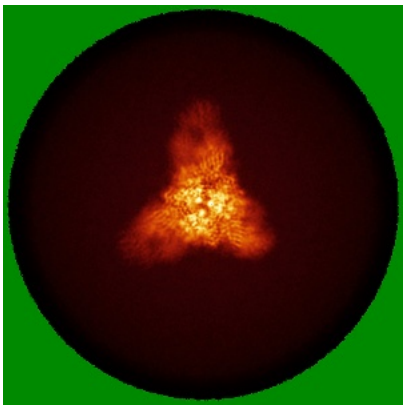
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

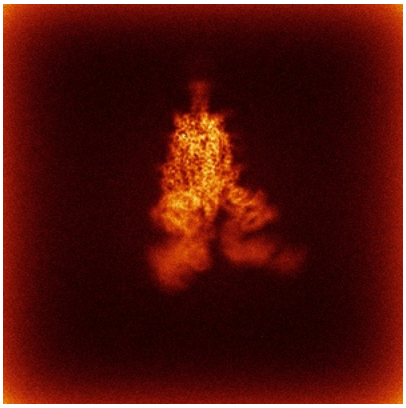
6.4.1 Primary map

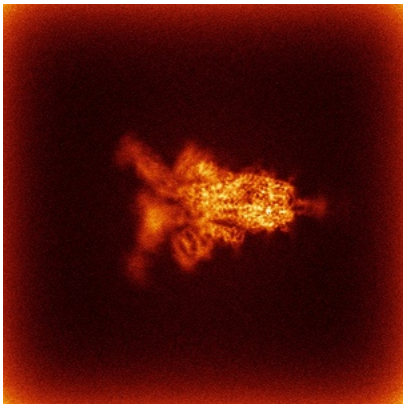

X

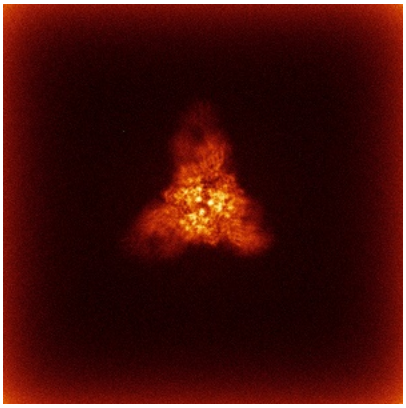

Y


Z

6.4.2 Raw map


X

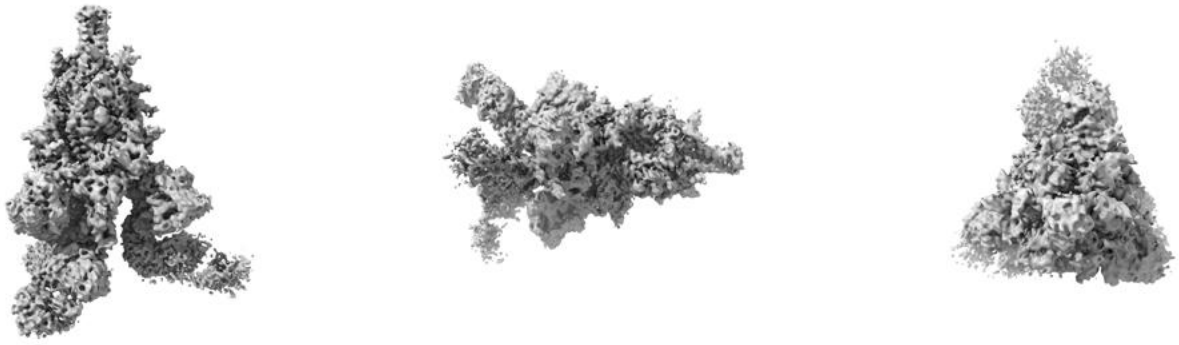

Y


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



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.2206. 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

X Y Z

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

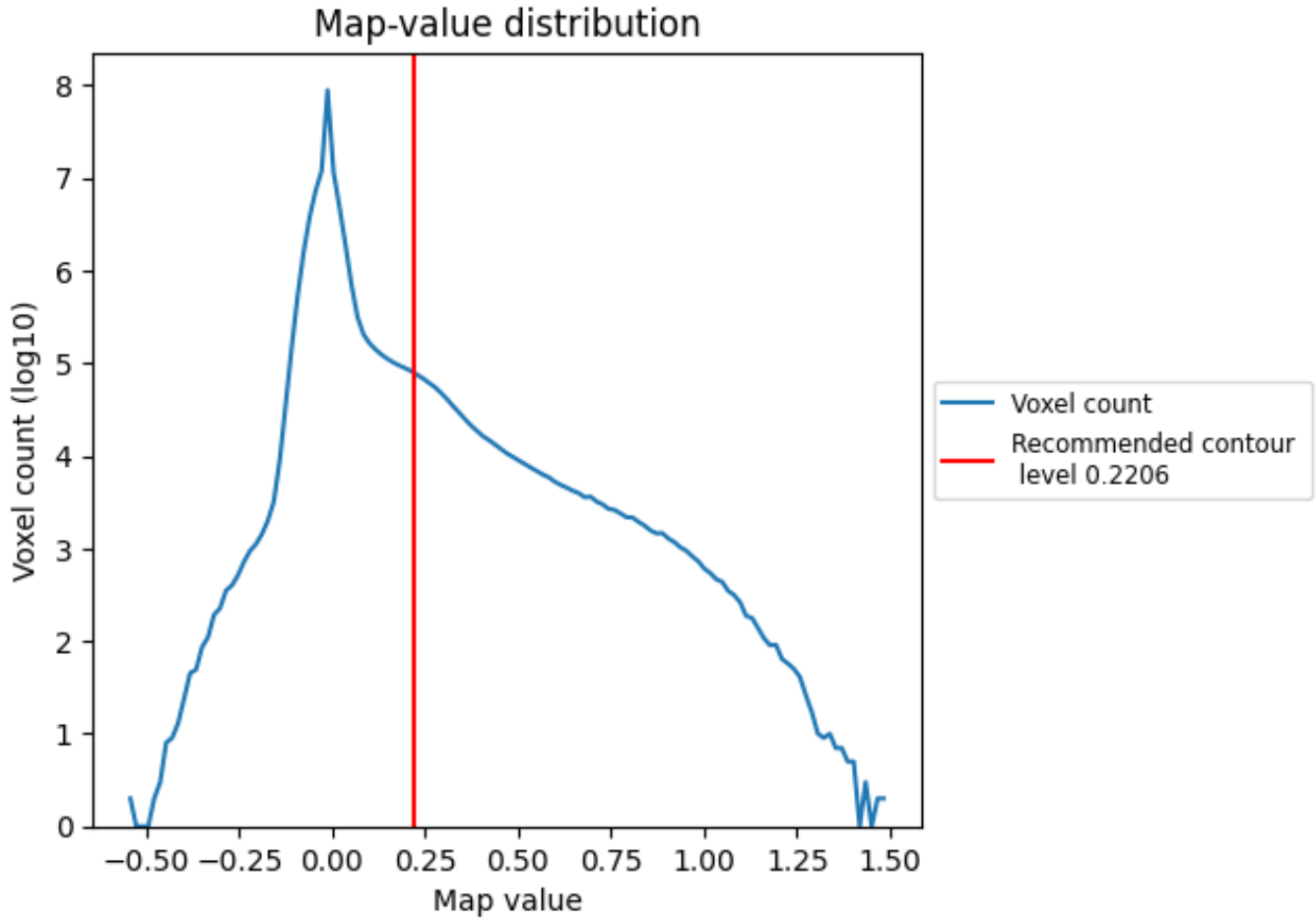
6.6 Mask visualisation (i)

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

7 Map analysis

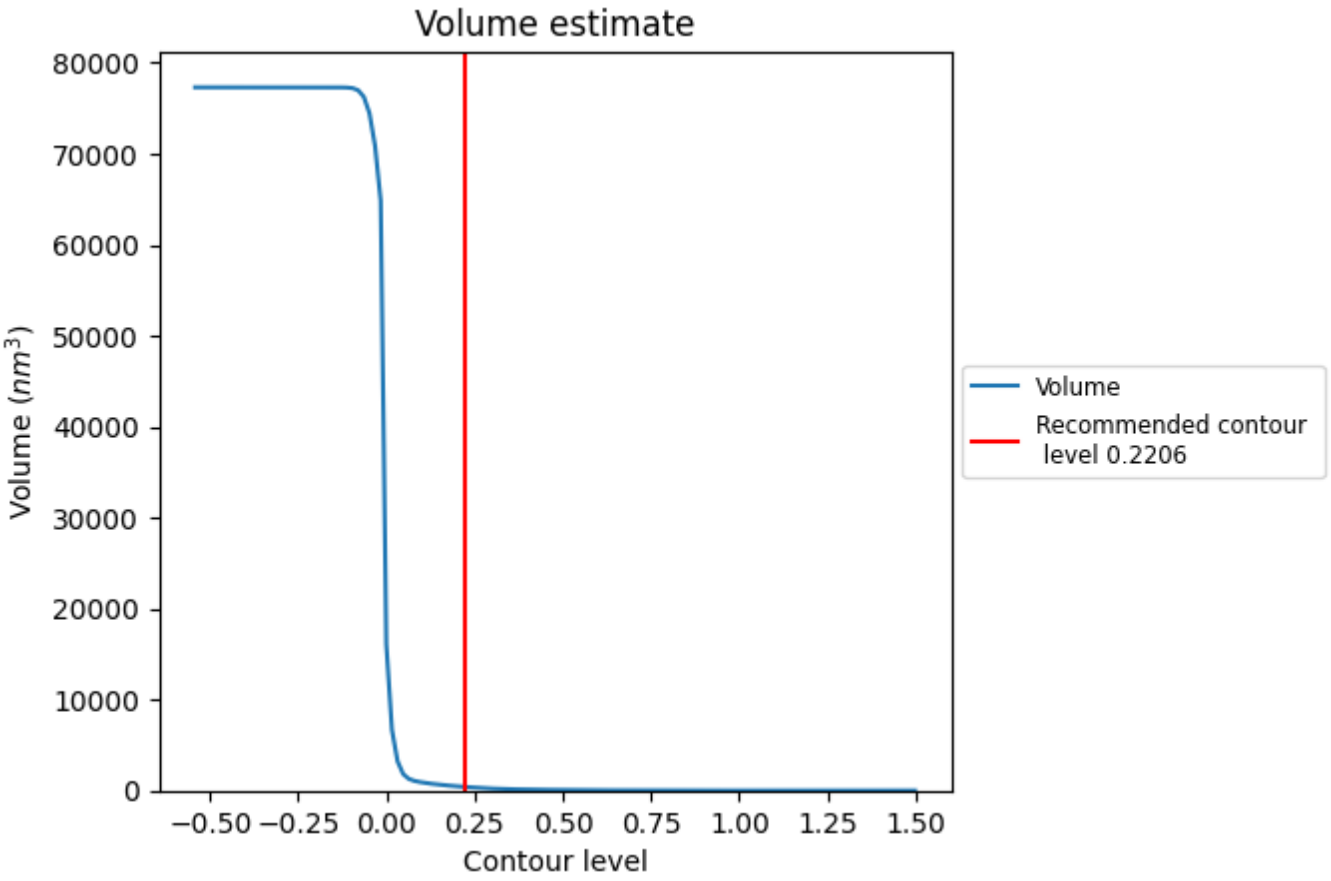
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

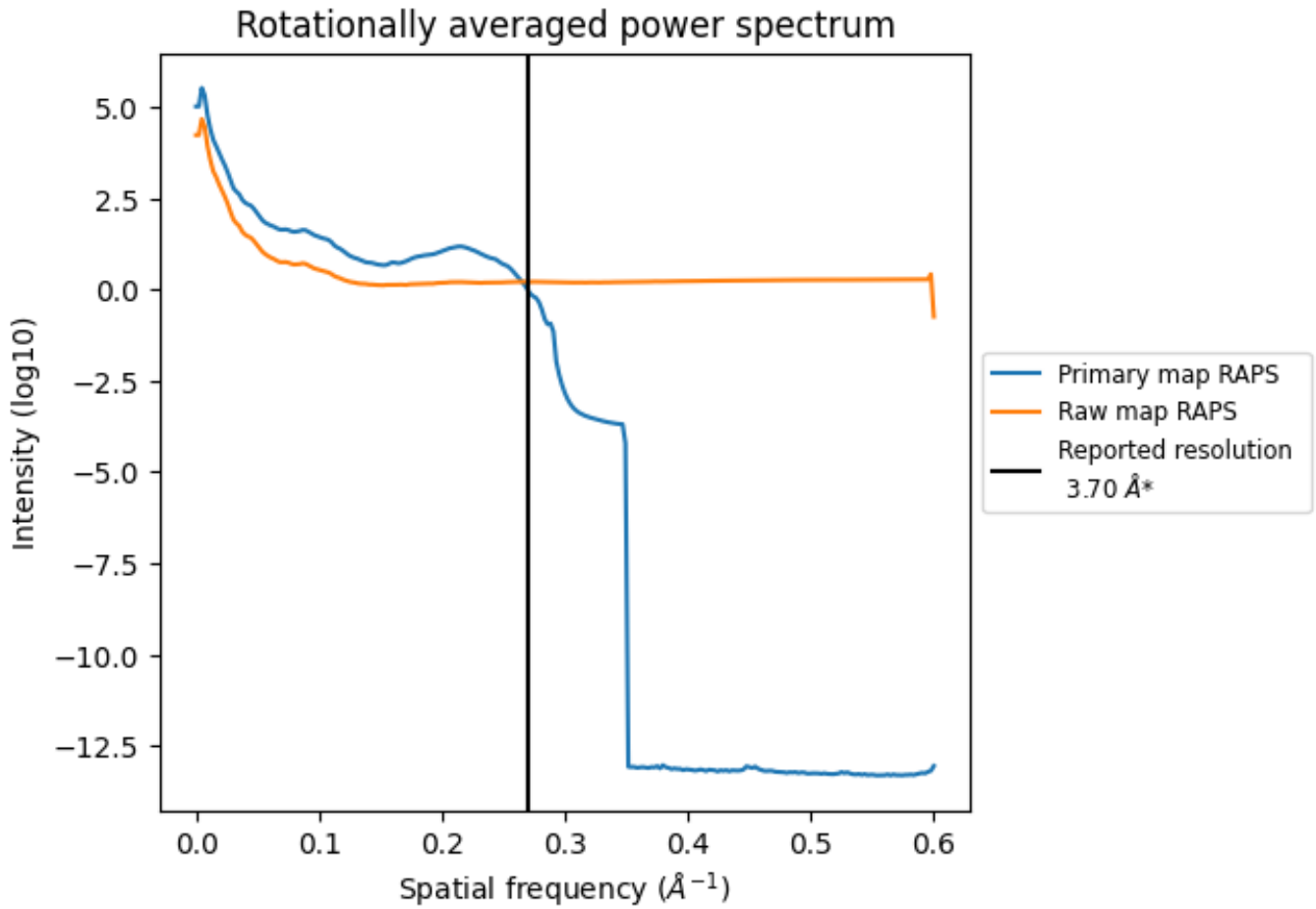
7.2 Volume estimate



The volume at the recommended contour level is 406 nm^3; this corresponds to an approximate mass of 366 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 (i)

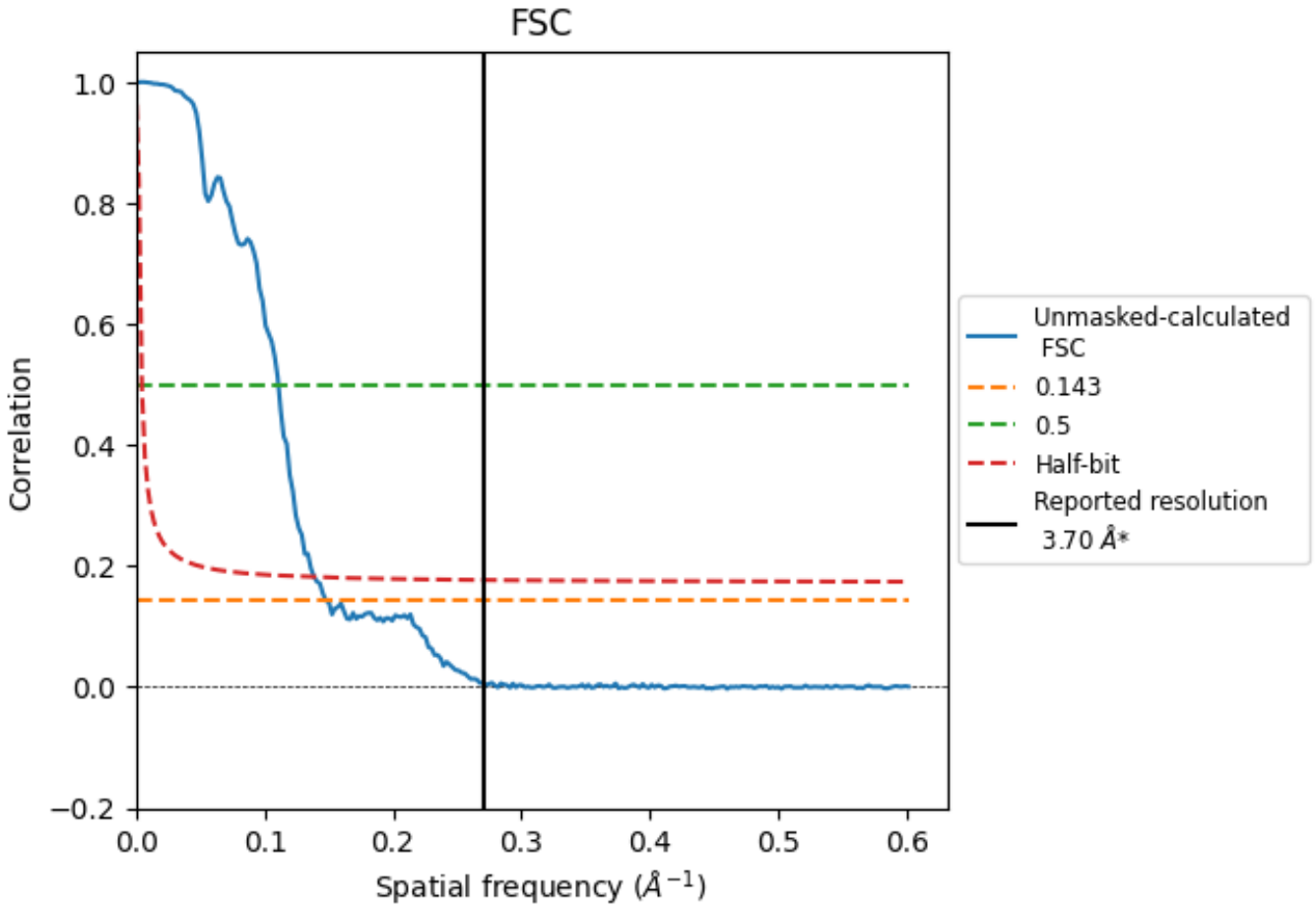


*Reported resolution corresponds to spatial frequency of 0.270 $Å^{-1}$

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.270 Å^{-1}

8.2 Resolution estimates (i)

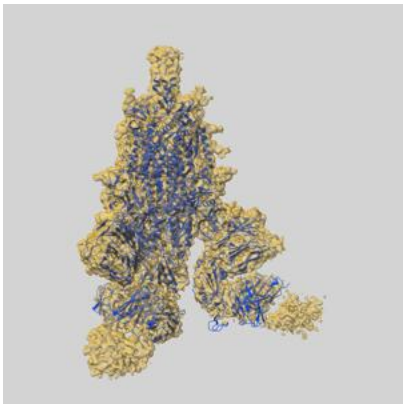
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	6.77	9.02	7.20

*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 6.77 differs from the reported value 3.7 by more than 10 %

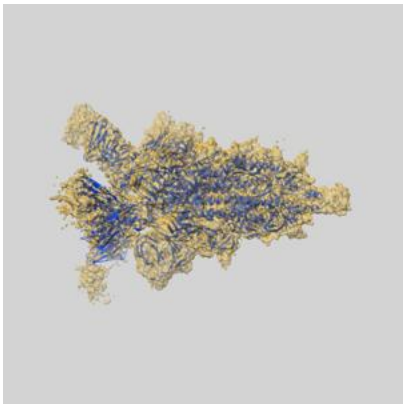
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-34408 and PDB model 8H01. Per-residue inclusion information can be found in section 3 on page 16.

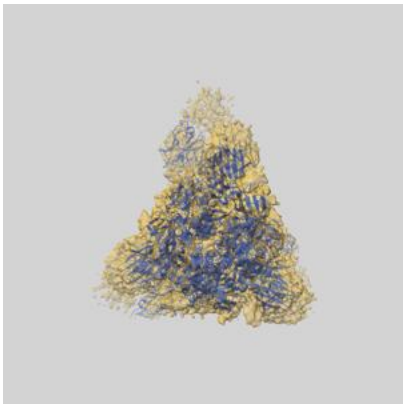
9.1 Map-model overlay



X



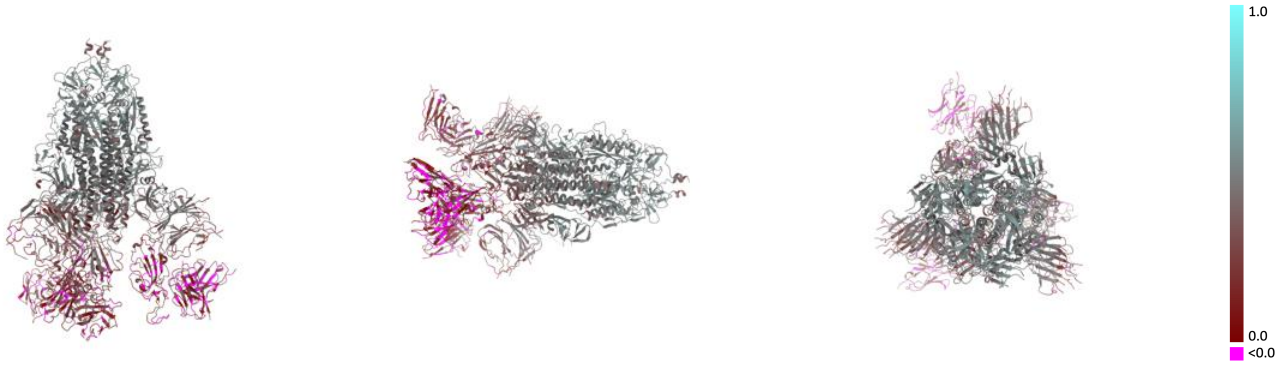
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.2206 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)



X Y Z

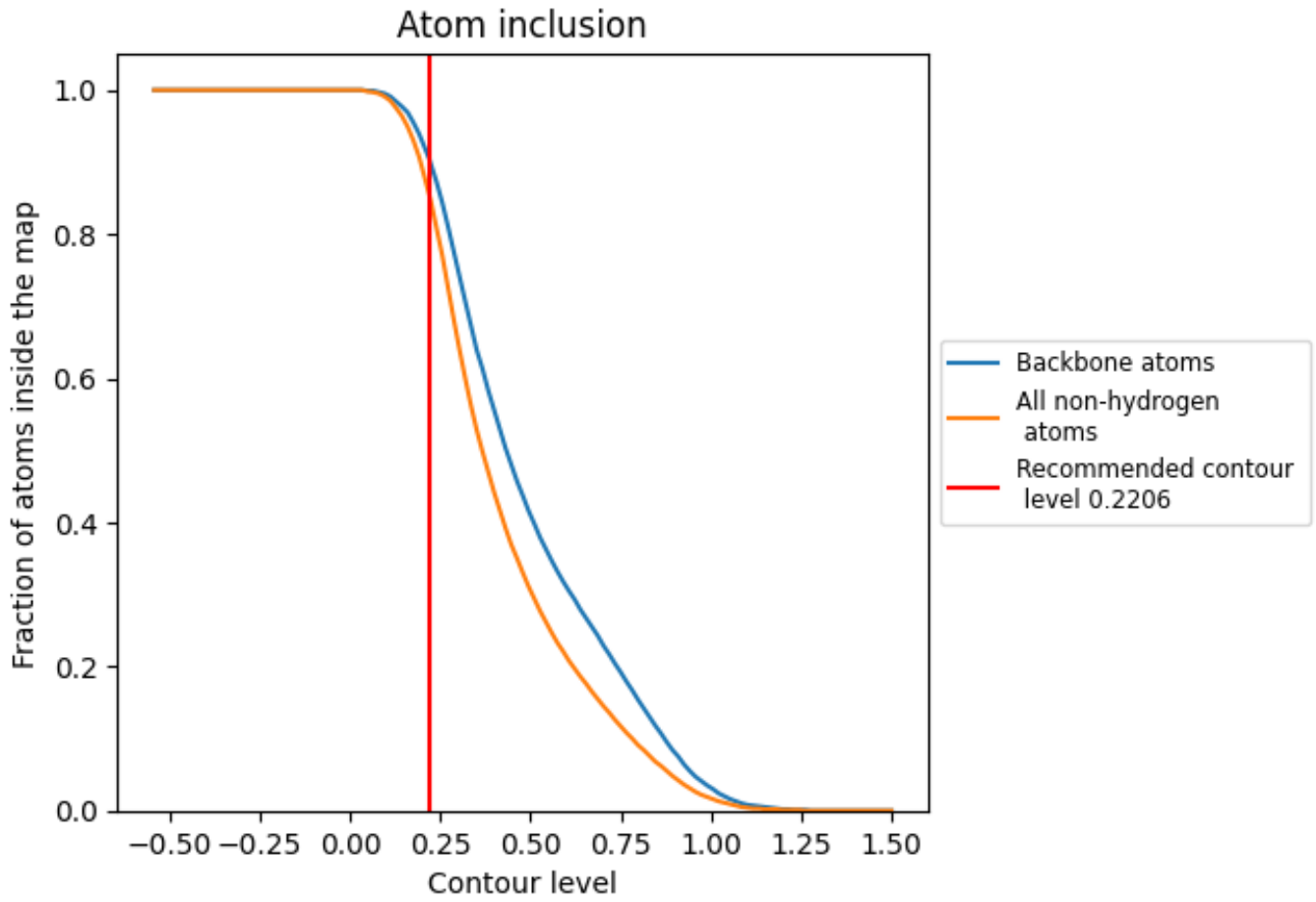
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

X Y Z

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 (0.2206).

9.4 Atom inclusion (i)



At the recommended contour level, 90% of all backbone atoms, 85% 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 (0.2206) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8530	0.3420
A	0.9170	0.4110
B	0.8820	0.3680
C	0.8900	0.3750
D	0.8240	0.2120
E	0.7990	0.2050
F	0.5300	0.0990
G	0.5810	0.0660
H	0.5970	0.1100
I	0.5560	0.1030
J	0.8930	0.4220
K	0.7140	0.4080
L	0.9640	0.3910
M	0.9290	0.4590
N	0.9640	0.4910
O	0.8570	0.4580
P	1.0000	0.4520
Q	0.6430	0.3060
R	0.8930	0.4010
S	0.8930	0.4170
T	0.9640	0.4940
U	0.8210	0.3780
V	0.5360	0.3470
W	1.0000	0.4260
X	0.8210	0.3800