



Full wwPDB EM Validation Report ⓘ

Aug 4, 2026 – 03:41 PM JST

PDB ID : 9XI0 / pdb_00009xi0
EMDB ID : EMD-66892
Title : SARS-CoV-2 Omicron BA.4/5 spike trimer in complex with C092 Fab and ACE2 (3 RBD up)
Authors : Niu, C.; Liu, B.; Gao, X.; Li, Z.; He, J.; Xiong, X.
Deposited on : 2025-11-02
Resolution : 3.98 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.dev133
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

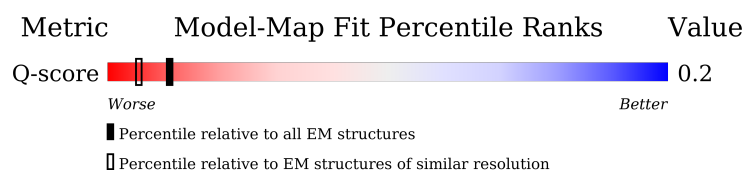
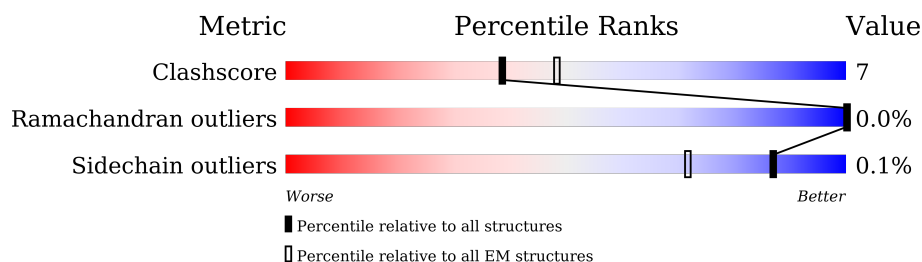
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.98 Å.

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



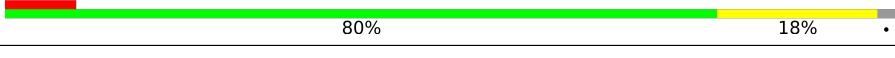
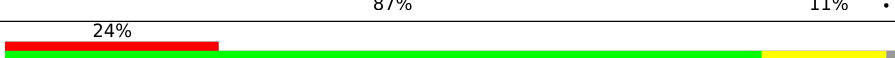
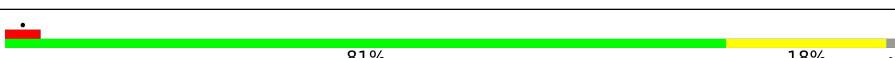
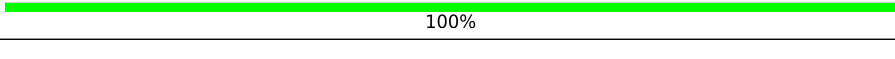
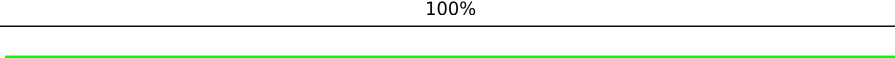
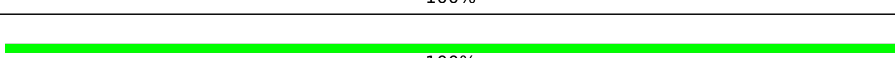
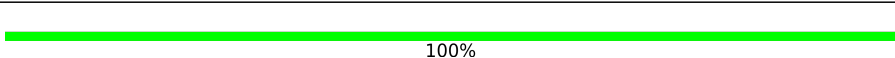
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	7512 (3.48 - 4.48)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1247	
1	B	1247	
1	C	1247	
2	G	228	

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Mol	Chain	Length	Quality of chain
2	H	228	
2	I	228	
3	J	217	
3	K	217	
3	L	217	
4	D	602	
4	E	602	
4	F	602	
5	M	2	
5	N	2	
5	O	2	
5	P	2	
5	Q	2	
5	R	2	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 49642 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	1051	Total	C	N	O	S	0	0
			8242	5271	1376	1558	37		
1	B	1051	Total	C	N	O	S	0	0
			8244	5273	1376	1558	37		
1	C	1051	Total	C	N	O	S	0	0
			8244	5273	1376	1558	37		

There are 249 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	ILE	THR	variant	UNP P0DTC2
A	27	SER	ALA	variant	UNP P0DTC2
A	142	ASP	GLY	variant	UNP P0DTC2
A	213	GLY	VAL	variant	UNP P0DTC2
A	339	ASP	GLY	variant	UNP P0DTC2
A	371	PHE	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	376	ALA	THR	variant	UNP P0DTC2
A	405	ASN	ASP	variant	UNP P0DTC2
A	408	SER	ARG	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	452	ARG	LEU	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	486	VAL	PHE	variant	UNP P0DTC2
A	498	ARG	GLN	variant	UNP P0DTC2
A	501	TYR	ASN	variant	UNP P0DTC2
A	505	HIS	TYR	variant	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	655	TYR	HIS	variant	UNP P0DTC2
A	679	LYS	ASN	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	817	PRO	PHE	engineered mutation	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	986	PRO	LYS	engineered mutation	UNP P0DTC2
A	987	PRO	VAL	engineered mutation	UNP P0DTC2
A	1212	GLY	-	expression tag	UNP P0DTC2
A	1213	SER	-	expression tag	UNP P0DTC2
A	1214	GLY	-	expression tag	UNP P0DTC2
A	1215	ARG	-	expression tag	UNP P0DTC2
A	1216	GLU	-	expression tag	UNP P0DTC2
A	1217	ASN	-	expression tag	UNP P0DTC2
A	1218	LEU	-	expression tag	UNP P0DTC2
A	1219	TYR	-	expression tag	UNP P0DTC2
A	1220	PHE	-	expression tag	UNP P0DTC2
A	1221	GLN	-	expression tag	UNP P0DTC2
A	1222	GLY	-	expression tag	UNP P0DTC2
A	1223	GLY	-	expression tag	UNP P0DTC2
A	1224	GLY	-	expression tag	UNP P0DTC2
A	1225	GLY	-	expression tag	UNP P0DTC2
A	1226	SER	-	expression tag	UNP P0DTC2
A	1227	GLY	-	expression tag	UNP P0DTC2
A	1228	TYR	-	expression tag	UNP P0DTC2
A	1229	ILE	-	expression tag	UNP P0DTC2
A	1230	PRO	-	expression tag	UNP P0DTC2
A	1231	GLU	-	expression tag	UNP P0DTC2
A	1232	ALA	-	expression tag	UNP P0DTC2
A	1233	PRO	-	expression tag	UNP P0DTC2
A	1234	ARG	-	expression tag	UNP P0DTC2
A	1235	ASP	-	expression tag	UNP P0DTC2
A	1236	GLY	-	expression tag	UNP P0DTC2
A	1237	GLN	-	expression tag	UNP P0DTC2
A	1238	ALA	-	expression tag	UNP P0DTC2
A	1239	TYR	-	expression tag	UNP P0DTC2
A	1240	VAL	-	expression tag	UNP P0DTC2
A	1241	ARG	-	expression tag	UNP P0DTC2
A	1242	LYS	-	expression tag	UNP P0DTC2
A	1243	ASP	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1244	GLY	-	expression tag	UNP P0DTC2
A	1245	GLU	-	expression tag	UNP P0DTC2
A	1246	TRP	-	expression tag	UNP P0DTC2
A	1247	VAL	-	expression tag	UNP P0DTC2
A	1248	LEU	-	expression tag	UNP P0DTC2
A	1249	LEU	-	expression tag	UNP P0DTC2
A	1250	SER	-	expression tag	UNP P0DTC2
A	1251	THR	-	expression tag	UNP P0DTC2
A	1252	PHE	-	expression tag	UNP P0DTC2
A	1253	LEU	-	expression tag	UNP P0DTC2
A	1254	GLY	-	expression tag	UNP P0DTC2
A	1255	HIS	-	expression tag	UNP P0DTC2
A	1256	HIS	-	expression tag	UNP P0DTC2
A	1257	HIS	-	expression tag	UNP P0DTC2
A	1258	HIS	-	expression tag	UNP P0DTC2
A	1259	HIS	-	expression tag	UNP P0DTC2
A	1260	HIS	-	expression tag	UNP P0DTC2
B	19	ILE	THR	variant	UNP P0DTC2
B	27	SER	ALA	variant	UNP P0DTC2
B	142	ASP	GLY	variant	UNP P0DTC2
B	213	GLY	VAL	variant	UNP P0DTC2
B	339	ASP	GLY	variant	UNP P0DTC2
B	371	PHE	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	376	ALA	THR	variant	UNP P0DTC2
B	405	ASN	ASP	variant	UNP P0DTC2
B	408	SER	ARG	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	452	ARG	LEU	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	486	VAL	PHE	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	505	HIS	TYR	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2
B	679	LYS	ASN	variant	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	796	TYR	ASP	variant	UNP P0DTC2
B	817	PRO	PHE	engineered mutation	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	986	PRO	LYS	engineered mutation	UNP P0DTC2
B	987	PRO	VAL	engineered mutation	UNP P0DTC2
B	1212	GLY	-	expression tag	UNP P0DTC2
B	1213	SER	-	expression tag	UNP P0DTC2
B	1214	GLY	-	expression tag	UNP P0DTC2
B	1215	ARG	-	expression tag	UNP P0DTC2
B	1216	GLU	-	expression tag	UNP P0DTC2
B	1217	ASN	-	expression tag	UNP P0DTC2
B	1218	LEU	-	expression tag	UNP P0DTC2
B	1219	TYR	-	expression tag	UNP P0DTC2
B	1220	PHE	-	expression tag	UNP P0DTC2
B	1221	GLN	-	expression tag	UNP P0DTC2
B	1222	GLY	-	expression tag	UNP P0DTC2
B	1223	GLY	-	expression tag	UNP P0DTC2
B	1224	GLY	-	expression tag	UNP P0DTC2
B	1225	GLY	-	expression tag	UNP P0DTC2
B	1226	SER	-	expression tag	UNP P0DTC2
B	1227	GLY	-	expression tag	UNP P0DTC2
B	1228	TYR	-	expression tag	UNP P0DTC2
B	1229	ILE	-	expression tag	UNP P0DTC2
B	1230	PRO	-	expression tag	UNP P0DTC2
B	1231	GLU	-	expression tag	UNP P0DTC2
B	1232	ALA	-	expression tag	UNP P0DTC2
B	1233	PRO	-	expression tag	UNP P0DTC2
B	1234	ARG	-	expression tag	UNP P0DTC2
B	1235	ASP	-	expression tag	UNP P0DTC2
B	1236	GLY	-	expression tag	UNP P0DTC2
B	1237	GLN	-	expression tag	UNP P0DTC2
B	1238	ALA	-	expression tag	UNP P0DTC2
B	1239	TYR	-	expression tag	UNP P0DTC2
B	1240	VAL	-	expression tag	UNP P0DTC2
B	1241	ARG	-	expression tag	UNP P0DTC2
B	1242	LYS	-	expression tag	UNP P0DTC2
B	1243	ASP	-	expression tag	UNP P0DTC2
B	1244	GLY	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1245	GLU	-	expression tag	UNP P0DTC2
B	1246	TRP	-	expression tag	UNP P0DTC2
B	1247	VAL	-	expression tag	UNP P0DTC2
B	1248	LEU	-	expression tag	UNP P0DTC2
B	1249	LEU	-	expression tag	UNP P0DTC2
B	1250	SER	-	expression tag	UNP P0DTC2
B	1251	THR	-	expression tag	UNP P0DTC2
B	1252	PHE	-	expression tag	UNP P0DTC2
B	1253	LEU	-	expression tag	UNP P0DTC2
B	1254	GLY	-	expression tag	UNP P0DTC2
B	1255	HIS	-	expression tag	UNP P0DTC2
B	1256	HIS	-	expression tag	UNP P0DTC2
B	1257	HIS	-	expression tag	UNP P0DTC2
B	1258	HIS	-	expression tag	UNP P0DTC2
B	1259	HIS	-	expression tag	UNP P0DTC2
B	1260	HIS	-	expression tag	UNP P0DTC2
C	19	ILE	THR	variant	UNP P0DTC2
C	27	SER	ALA	variant	UNP P0DTC2
C	142	ASP	GLY	variant	UNP P0DTC2
C	213	GLY	VAL	variant	UNP P0DTC2
C	339	ASP	GLY	variant	UNP P0DTC2
C	371	PHE	SER	variant	UNP P0DTC2
C	373	PRO	SER	variant	UNP P0DTC2
C	375	PHE	SER	variant	UNP P0DTC2
C	376	ALA	THR	variant	UNP P0DTC2
C	405	ASN	ASP	variant	UNP P0DTC2
C	408	SER	ARG	variant	UNP P0DTC2
C	417	ASN	LYS	variant	UNP P0DTC2
C	440	LYS	ASN	variant	UNP P0DTC2
C	452	ARG	LEU	variant	UNP P0DTC2
C	477	ASN	SER	variant	UNP P0DTC2
C	478	LYS	THR	variant	UNP P0DTC2
C	484	ALA	GLU	variant	UNP P0DTC2
C	486	VAL	PHE	variant	UNP P0DTC2
C	498	ARG	GLN	variant	UNP P0DTC2
C	501	TYR	ASN	variant	UNP P0DTC2
C	505	HIS	TYR	variant	UNP P0DTC2
C	614	GLY	ASP	variant	UNP P0DTC2
C	655	TYR	HIS	variant	UNP P0DTC2
C	679	LYS	ASN	variant	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	817	PRO	PHE	engineered mutation	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	986	PRO	LYS	engineered mutation	UNP P0DTC2
C	987	PRO	VAL	engineered mutation	UNP P0DTC2
C	1212	GLY	-	expression tag	UNP P0DTC2
C	1213	SER	-	expression tag	UNP P0DTC2
C	1214	GLY	-	expression tag	UNP P0DTC2
C	1215	ARG	-	expression tag	UNP P0DTC2
C	1216	GLU	-	expression tag	UNP P0DTC2
C	1217	ASN	-	expression tag	UNP P0DTC2
C	1218	LEU	-	expression tag	UNP P0DTC2
C	1219	TYR	-	expression tag	UNP P0DTC2
C	1220	PHE	-	expression tag	UNP P0DTC2
C	1221	GLN	-	expression tag	UNP P0DTC2
C	1222	GLY	-	expression tag	UNP P0DTC2
C	1223	GLY	-	expression tag	UNP P0DTC2
C	1224	GLY	-	expression tag	UNP P0DTC2
C	1225	GLY	-	expression tag	UNP P0DTC2
C	1226	SER	-	expression tag	UNP P0DTC2
C	1227	GLY	-	expression tag	UNP P0DTC2
C	1228	TYR	-	expression tag	UNP P0DTC2
C	1229	ILE	-	expression tag	UNP P0DTC2
C	1230	PRO	-	expression tag	UNP P0DTC2
C	1231	GLU	-	expression tag	UNP P0DTC2
C	1232	ALA	-	expression tag	UNP P0DTC2
C	1233	PRO	-	expression tag	UNP P0DTC2
C	1234	ARG	-	expression tag	UNP P0DTC2
C	1235	ASP	-	expression tag	UNP P0DTC2
C	1236	GLY	-	expression tag	UNP P0DTC2
C	1237	GLN	-	expression tag	UNP P0DTC2
C	1238	ALA	-	expression tag	UNP P0DTC2
C	1239	TYR	-	expression tag	UNP P0DTC2
C	1240	VAL	-	expression tag	UNP P0DTC2
C	1241	ARG	-	expression tag	UNP P0DTC2
C	1242	LYS	-	expression tag	UNP P0DTC2
C	1243	ASP	-	expression tag	UNP P0DTC2
C	1244	GLY	-	expression tag	UNP P0DTC2
C	1245	GLU	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1246	TRP	-	expression tag	UNP P0DTC2
C	1247	VAL	-	expression tag	UNP P0DTC2
C	1248	LEU	-	expression tag	UNP P0DTC2
C	1249	LEU	-	expression tag	UNP P0DTC2
C	1250	SER	-	expression tag	UNP P0DTC2
C	1251	THR	-	expression tag	UNP P0DTC2
C	1252	PHE	-	expression tag	UNP P0DTC2
C	1253	LEU	-	expression tag	UNP P0DTC2
C	1254	GLY	-	expression tag	UNP P0DTC2
C	1255	HIS	-	expression tag	UNP P0DTC2
C	1256	HIS	-	expression tag	UNP P0DTC2
C	1257	HIS	-	expression tag	UNP P0DTC2
C	1258	HIS	-	expression tag	UNP P0DTC2
C	1259	HIS	-	expression tag	UNP P0DTC2
C	1260	HIS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called Heavy chain of C092 Fab.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	222	Total	C	N	O	S	2	0
			1645	1035	275	326	9		
2	G	222	Total	C	N	O	S	2	0
			1645	1035	275	326	9		
2	I	222	Total	C	N	O	S	2	0
			1645	1035	275	326	9		

- Molecule 3 is a protein called Light chain of C092 Fab.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	L	213	Total	C	N	O	S	0	0
			1570	978	263	325	4		
3	J	213	Total	C	N	O	S	0	0
			1570	978	263	325	4		
3	K	213	Total	C	N	O	S	0	0
			1570	978	263	325	4		

- Molecule 4 is a protein called Processed angiotensin-converting enzyme 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	596	Total	C	N	O	S	0	0
			4864	3112	805	918	29		

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Mol	Chain	Residues	Atoms					AltConf	Trace
4	F	596	Total	C	N	O	S	0	0
			4864	3112	805	918	29		
4	D	596	Total	C	N	O	S	0	0
			4864	3112	805	918	29		

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	616	PRO	-	expression tag	UNP Q9BYF1
E	617	LEU	-	expression tag	UNP Q9BYF1
E	618	VAL	-	expression tag	UNP Q9BYF1
E	619	PRO	-	expression tag	UNP Q9BYF1
E	620	ARG	-	expression tag	UNP Q9BYF1
F	616	PRO	-	expression tag	UNP Q9BYF1
F	617	LEU	-	expression tag	UNP Q9BYF1
F	618	VAL	-	expression tag	UNP Q9BYF1
F	619	PRO	-	expression tag	UNP Q9BYF1
F	620	ARG	-	expression tag	UNP Q9BYF1
D	616	PRO	-	expression tag	UNP Q9BYF1
D	617	LEU	-	expression tag	UNP Q9BYF1
D	618	VAL	-	expression tag	UNP Q9BYF1
D	619	PRO	-	expression tag	UNP Q9BYF1
D	620	ARG	-	expression tag	UNP Q9BYF1

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



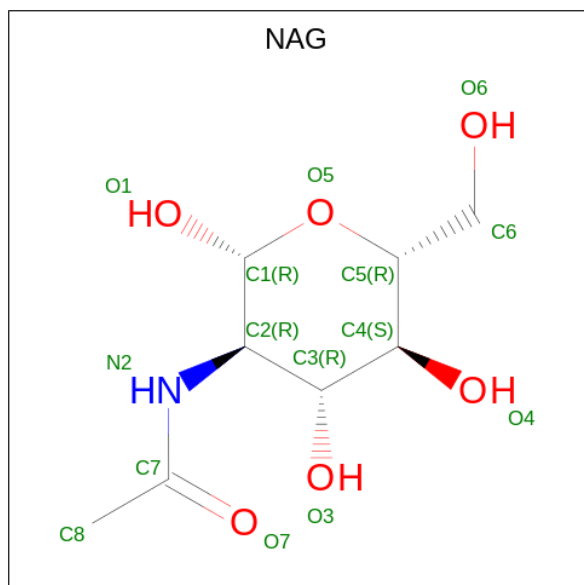
Mol	Chain	Residues	Atoms				AltConf	Trace
5	M	2	Total	C	N	O	0	0
			28	16	2	10		
5	N	2	Total	C	N	O	0	0
			28	16	2	10		
5	O	2	Total	C	N	O	0	0
			28	16	2	10		
5	P	2	Total	C	N	O	0	0
			28	16	2	10		
5	Q	2	Total	C	N	O	0	0
			28	16	2	10		

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

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

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Mol	Chain	Residues	Atoms				AltConf
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	B	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	C	1	Total	C	N	O	0
			14	8	1	5	
6	E	1	Total	C	N	O	0
			14	8	1	5	
6	E	1	Total	C	N	O	0
			14	8	1	5	
6	E	1	Total	C	N	O	0
			14	8	1	5	
6	E	1	Total	C	N	O	0
			14	8	1	5	
6	F	1	Total	C	N	O	0
			14	8	1	5	
6	F	1	Total	C	N	O	0
			14	8	1	5	
6	F	1	Total	C	N	O	0
			14	8	1	5	

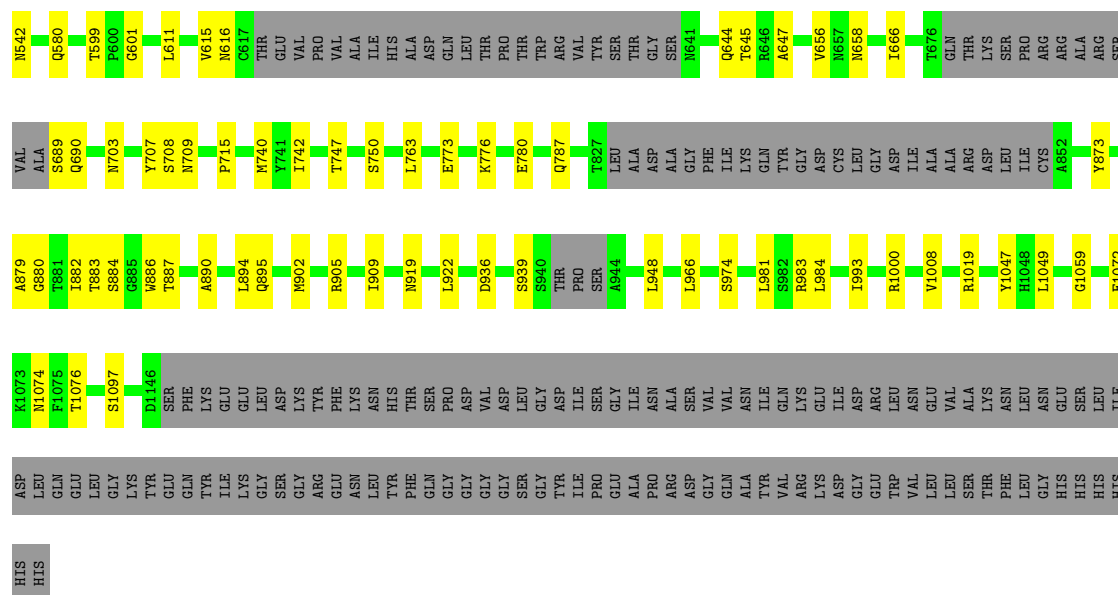
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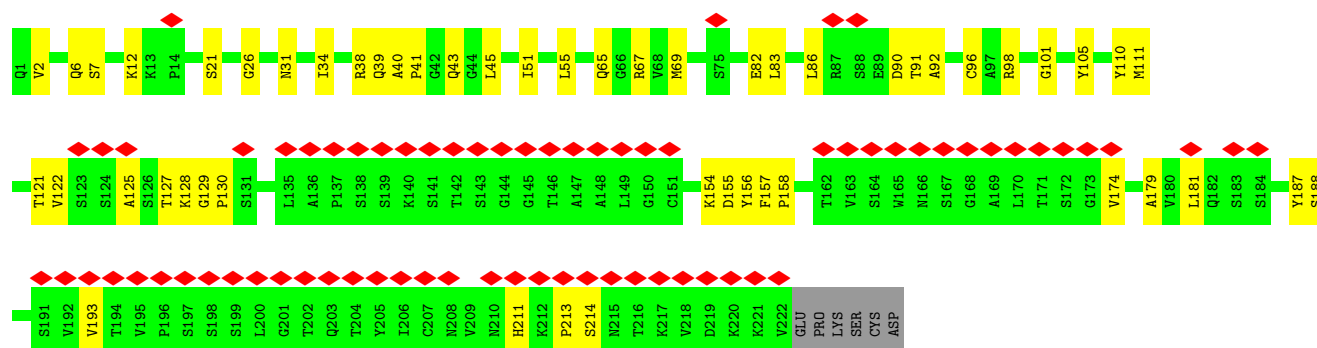
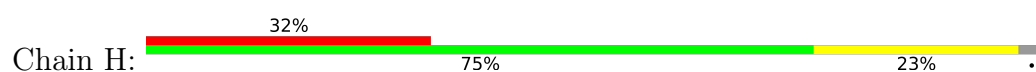
Mol	Chain	Residues	Atoms				AltConf
6	F	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

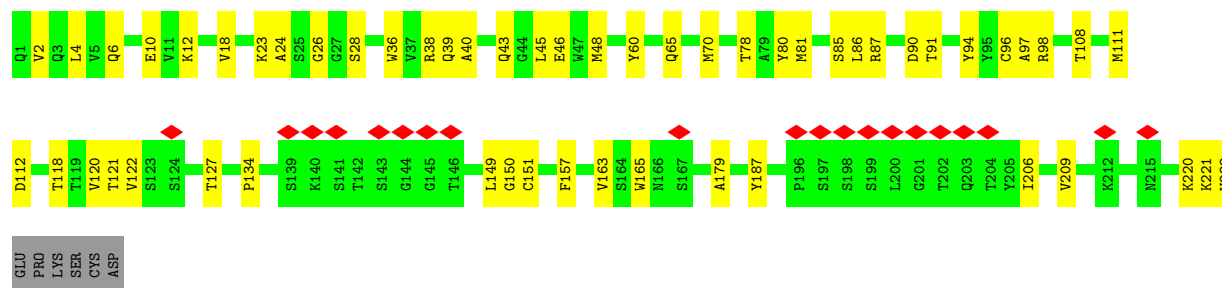
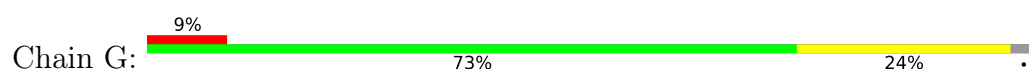
Mol	Chain	Residues	Atoms		AltConf
7	E	1	Total	Zn	0
			1	1	
7	F	1	Total	Zn	0
			1	1	
7	D	1	Total	Zn	0
			1	1	



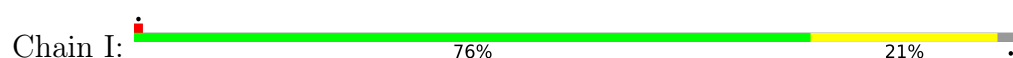
• Molecule 2: Heavy chain of C092 Fab

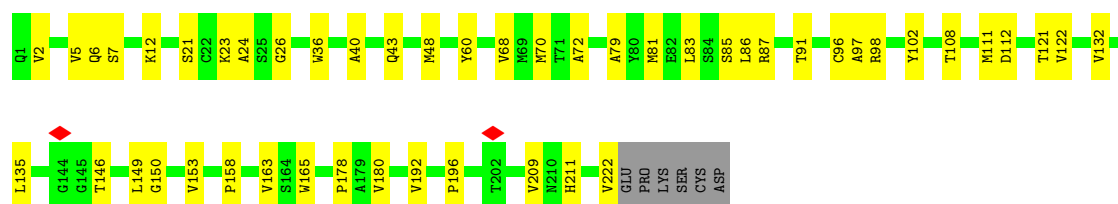


• Molecule 2: Heavy chain of C092 Fab

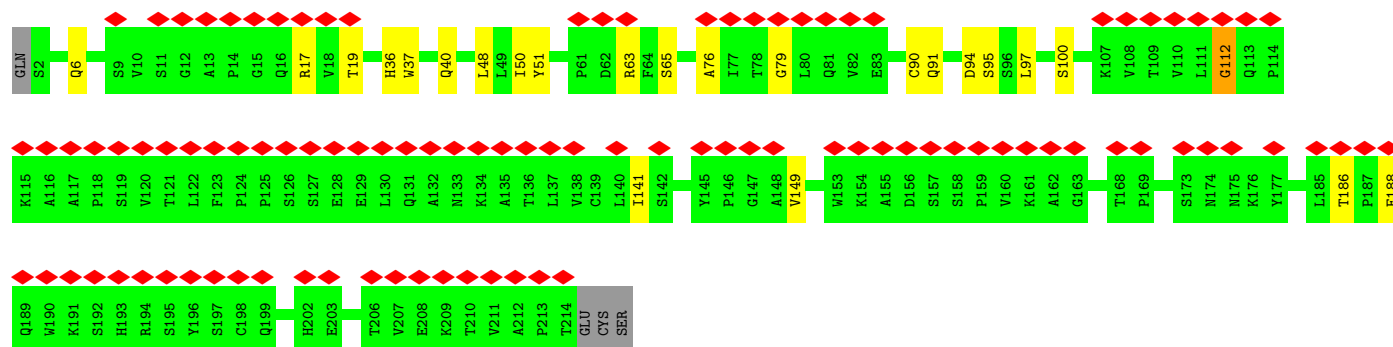
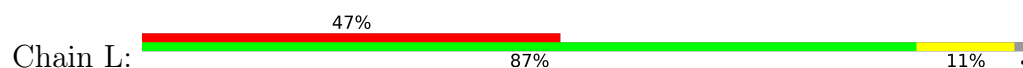


• Molecule 2: Heavy chain of C092 Fab

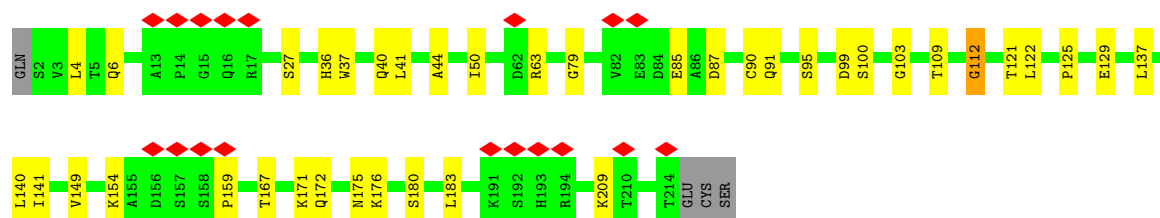




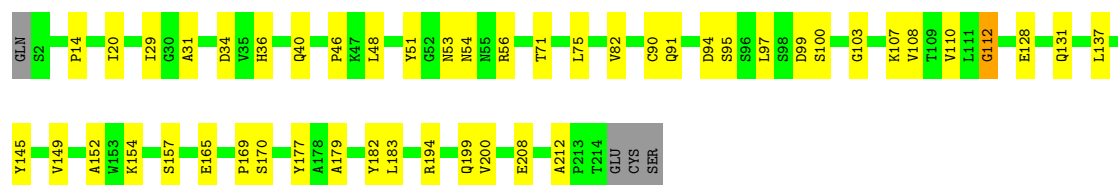
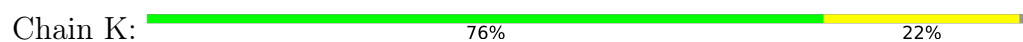
• Molecule 3: Light chain of C092 Fab



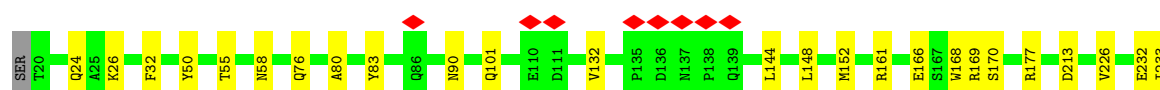
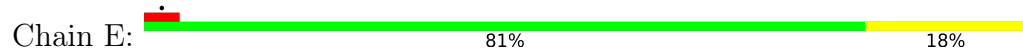
• Molecule 3: Light chain of C092 Fab

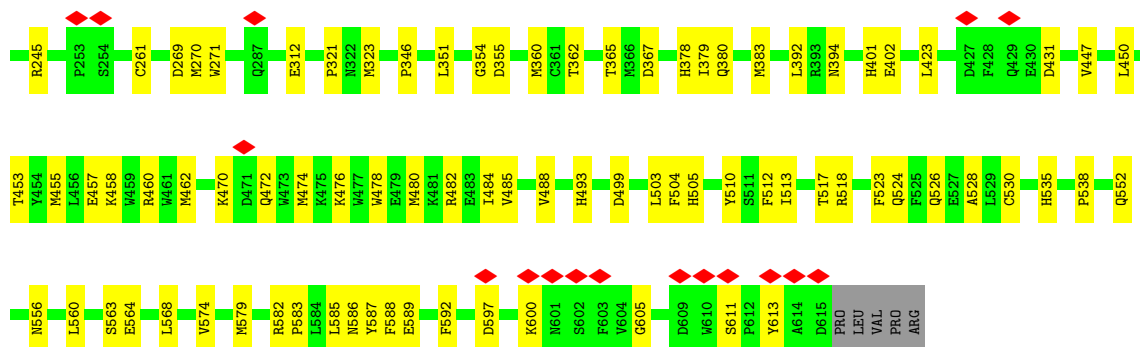


• Molecule 3: Light chain of C092 Fab



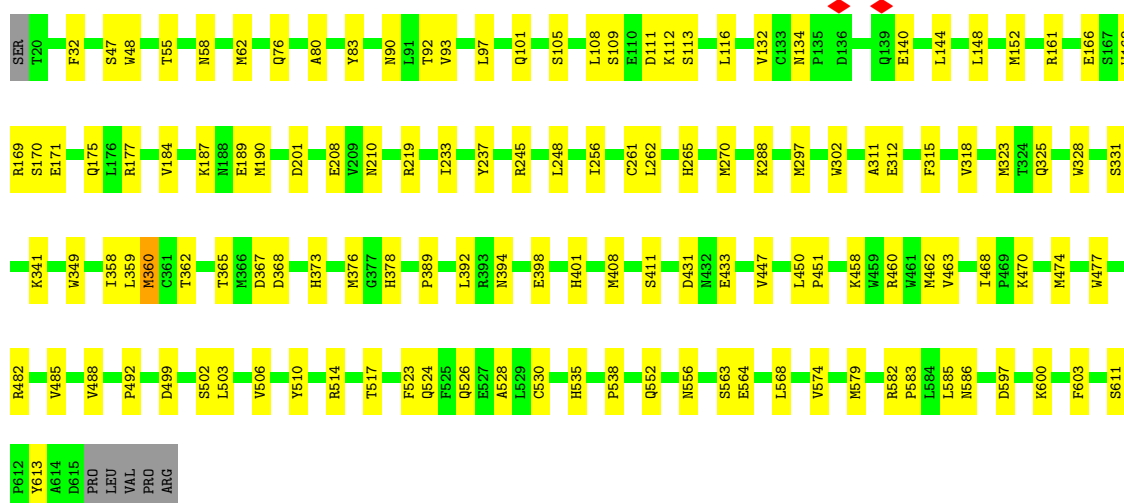
• Molecule 4: Processed angiotensin-converting enzyme 2





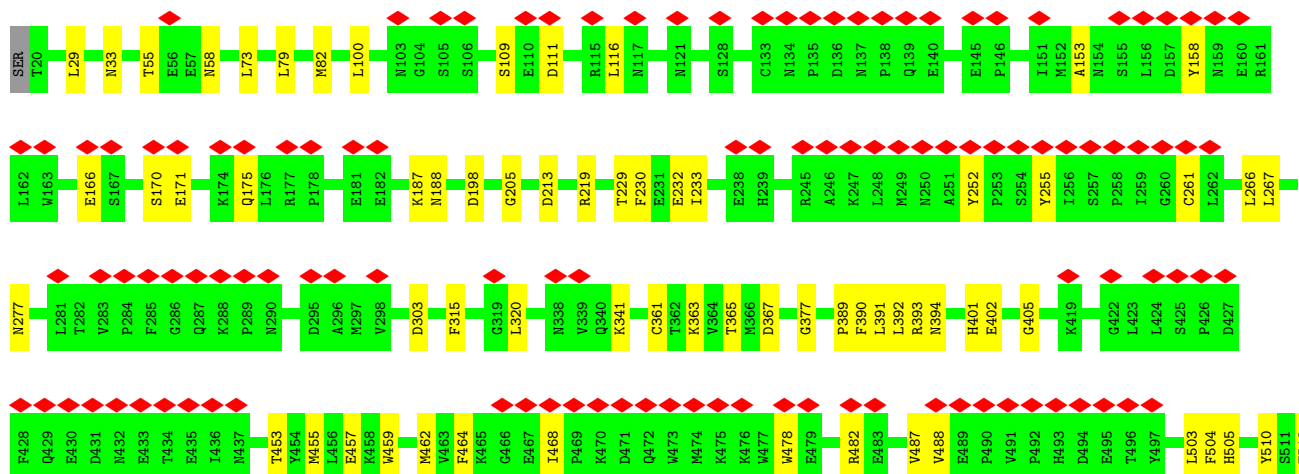
• Molecule 4: Processed angiotensin-converting enzyme 2

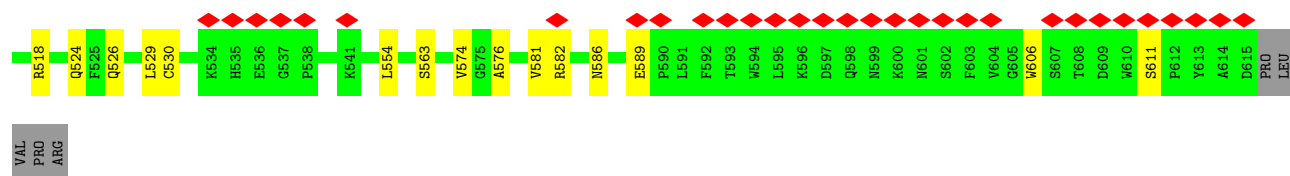
Chain F: 78% 21% .



• Molecule 4: Processed angiotensin-converting enzyme 2

Chain D: 24% 85% 14% .





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

Chain M:  50%



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

Chain N:  100%



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

Chain O:  100%



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

Chain P:  100%



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

Chain Q:  100%



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

Chain R:  100%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	148642	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.105	Depositor
Minimum map value	-0.366	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.029	Depositor
Recommended contour level	0.07	Depositor
Map size (Å)	422.4, 422.4, 422.4	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.65, 1.65, 1.65	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, ZN

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	0.25	0/8439	0.43	1/11479 (0.0%)
1	B	0.25	0/8442	0.41	0/11484
1	C	0.22	0/8442	0.39	0/11484
2	G	0.20	0/1691	0.43	0/2301
2	H	0.23	0/1691	0.44	0/2301
2	I	0.18	0/1691	0.40	0/2301
3	J	0.15	0/1608	0.49	1/2201 (0.0%)
3	K	0.13	0/1608	0.51	1/2201 (0.0%)
3	L	0.13	0/1608	0.48	1/2201 (0.0%)
4	D	0.13	0/5001	0.32	0/6795
4	E	0.15	0/5001	0.37	0/6795
4	F	0.20	0/5001	0.41	0/6795
All	All	0.21	0/50223	0.41	4/68338 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
3	J	0	1
3	K	0	1
3	L	0	1
All	All	0	3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	112	GLY	O-C-N	-17.18	100.37	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	112	GLY	O-C-N	-16.44	101.33	122.70
3	J	112	GLY	O-C-N	-14.73	103.55	122.70
1	A	527	PRO	N-CA-CB	6.33	109.89	103.25

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	J	112	GLY	Mainchain
3	K	112	GLY	Mainchain
3	L	112	GLY	Mainchain

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	8242	0	8033	123	0
1	B	8244	0	8039	128	0
1	C	8244	0	8039	118	0
2	G	1645	0	1614	37	0
2	H	1645	0	1614	35	0
2	I	1645	0	1614	31	0
3	J	1570	0	1504	24	0
3	K	1570	0	1504	31	0
3	L	1570	0	1504	14	0
4	D	4864	0	4634	49	0
4	E	4864	0	4634	73	0
4	F	4864	0	4634	82	0
5	M	28	0	25	1	0
5	N	28	0	25	0	0
5	O	28	0	25	0	0
5	P	28	0	25	0	0
5	Q	28	0	25	0	0
5	R	28	0	25	0	0
6	A	112	0	104	0	0
6	B	112	0	104	2	0
6	C	112	0	104	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	D	56	0	52	0	0
6	E	56	0	52	0	0
6	F	56	0	52	0	0
7	D	1	0	0	0	0
7	E	1	0	0	0	0
7	F	1	0	0	0	0
All	All	49642	0	47985	710	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (710) 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)
4:F:362:THR:HG23	4:F:368:ASP:HB3	1.51	0.91
1:A:360:ASN:H	1:A:523:THR:HB	1.39	0.85
1:A:335:LEU:HD22	1:A:364:ASP:HB2	1.58	0.84
2:G:97:ALA:HB3	2:G:111:MET:HE3	1.60	0.84
1:C:355:ARG:HD2	1:C:396:TYR:HD1	1.45	0.81
1:C:329:PHE:HB2	1:C:330:PRO:HD3	1.62	0.80
1:C:355:ARG:HD2	1:C:396:TYR:CD1	2.17	0.79
2:I:97:ALA:HB3	2:I:111:MET:HE3	1.69	0.75
4:F:311:ALA:HB1	4:F:376:MET:HE1	1.68	0.74
1:B:561:PRO:HA	1:B:577:ARG:HH12	1.51	0.73
1:B:334:ASN:HB2	1:B:362:VAL:HB	1.71	0.72
1:B:332:ILE:HG13	6:B:1302:NAG:H82	1.69	0.72
2:H:154:LYS:HD2	2:H:188:SER:HB3	1.73	0.71
1:B:543:PHE:O	1:B:544:ASN:OD1	2.07	0.71
1:B:331:ASN:HA	1:B:580:GLN:HG3	1.74	0.70
1:B:327:VAL:HG22	1:B:542:ASN:HB3	1.74	0.69
1:B:329:PHE:O	1:B:579:PRO:HB2	1.93	0.69
4:D:261:CYS:HB2	4:D:488:VAL:HG23	1.75	0.68
4:F:474:MET:HA	4:F:477:TRP:HB3	1.77	0.67
1:C:331:ASN:HB3	1:C:580:GLN:HA	1.76	0.67
1:A:64:TRP:HE1	1:A:264:ALA:HB1	1.57	0.67
4:F:458:LYS:HG2	4:F:462:MET:HE1	1.75	0.66
1:B:1052:PHE:HB2	1:B:1063:LEU:HB2	1.78	0.66
1:B:31:SER:HA	1:B:216:LEU:HD12	1.76	0.66
4:D:462:MET:HG2	4:D:468:ILE:HG13	1.77	0.66
1:B:332:ILE:HA	6:B:1302:NAG:C7	2.26	0.66
1:C:909:ILE:HD12	1:C:1047:TYR:HB3	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:461:LEU:HD21	1:B:467:ASP:HB2	1.77	0.65
1:A:398:ASP:HB2	1:A:512:VAL:HB	1.78	0.65
1:B:64:TRP:HE1	1:B:264:ALA:HB1	1.61	0.65
3:J:122:LEU:H	3:J:209:LYS:HE3	1.62	0.65
4:F:373:HIS:HA	4:F:376:MET:HE2	1.77	0.65
4:D:377:GLY:HA3	4:D:405:GLY:HA2	1.79	0.64
1:A:134:GLN:HB2	1:A:161:SER:HB2	1.79	0.64
1:A:534:VAL:HG21	1:A:539:VAL:HG11	1.80	0.64
3:J:37:TRP:H	3:J:50:ILE:HB	1.62	0.64
2:H:2:VAL:HA	2:H:26:GLY:HA3	1.80	0.64
1:B:544:ASN:HD22	1:B:579:PRO:HB3	1.63	0.64
1:C:64:TRP:HE1	1:C:264:ALA:HB1	1.62	0.64
1:C:393:THR:HA	1:C:522:ALA:HA	1.80	0.63
2:G:6:GLN:NE2	2:G:96:CYS:SG	2.72	0.63
1:A:856:ASN:HD21	1:A:963:VAL:HG12	1.64	0.63
3:K:165:GLU:HB2	3:K:182:TYR:HB2	1.80	0.63
1:C:201:PHE:HB3	1:C:229:LEU:HB2	1.82	0.62
4:D:478:TRP:HB3	4:D:482:ARG:HH12	1.63	0.62
1:C:689:SER:OG	1:C:690:GLN:N	2.32	0.62
4:E:152:MET:O	4:E:161:ARG:NH1	2.33	0.61
4:F:168:TRP:HE1	4:F:502:SER:HB2	1.63	0.61
4:E:453:THR:HG23	4:E:512:PHE:HB3	1.82	0.61
2:I:40:ALA:HB3	2:I:43:GLN:HB2	1.81	0.61
1:A:922:LEU:HD11	5:M:1:NAG:H3	1.82	0.61
2:H:40:ALA:HB3	2:H:43:GLN:HB2	1.82	0.61
4:E:517:THR:HG22	4:E:579:MET:HE1	1.83	0.61
1:A:983:ARG:HG3	1:A:984:LEU:HG	1.83	0.61
1:A:201:PHE:HB3	1:A:229:LEU:HB2	1.83	0.60
1:A:328:ARG:HH12	1:A:533:LEU:HB2	1.65	0.60
1:C:152:TRP:HB2	1:C:179:LEU:HB3	1.81	0.60
4:F:184:VAL:HG21	4:F:463:VAL:HG12	1.83	0.60
1:A:106:PHE:HB2	1:A:117:LEU:HB2	1.82	0.60
1:C:66:HIS:HB2	1:C:78:ARG:HG2	1.84	0.60
1:C:246:ARG:HH22	1:C:254:SER:HB3	1.66	0.60
2:H:6:GLN:NE2	2:H:96:CYS:SG	2.74	0.60
1:A:97:LYS:HD3	1:A:187:LYS:H	1.66	0.60
4:E:394:ASN:OD1	4:E:401:HIS:NE2	2.34	0.60
4:E:523:PHE:HB3	4:E:583:PRO:HB2	1.84	0.60
1:B:333:THR:HG23	1:B:527:PRO:HB3	1.84	0.60
4:E:152:MET:HE1	4:E:270:MET:HE2	1.83	0.60
4:E:360:MET:SD	4:E:362:THR:OG1	2.60	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:7:SER:HB3	2:I:21:SER:H	1.67	0.59
2:G:40:ALA:HB3	2:G:43:GLN:HB2	1.83	0.59
4:F:169:ARG:HE	4:F:499:ASP:HB3	1.67	0.59
1:C:130:VAL:HB	1:C:168:PHE:HB3	1.85	0.59
4:E:458:LYS:HG2	4:E:462:MET:HE1	1.84	0.59
4:D:392:LEU:HD13	4:D:563:SER:HB3	1.85	0.59
1:C:329:PHE:CE2	1:C:528:LYS:HD3	2.37	0.58
2:I:2:VAL:HA	2:I:26:GLY:HA3	1.84	0.58
1:B:543:PHE:HD1	1:B:579:PRO:HD3	1.68	0.58
2:I:6:GLN:NE2	2:I:96:CYS:SG	2.77	0.58
1:B:543:PHE:O	1:B:544:ASN:CG	2.45	0.58
1:B:599:THR:HG22	1:B:601:GLY:H	1.67	0.58
1:B:656:VAL:HG22	1:B:658:ASN:H	1.67	0.58
1:C:214:ARG:NH1	1:C:266:TYR:OH	2.37	0.58
3:L:91:GLN:NE2	3:L:100:SER:OG	2.36	0.58
2:I:85:SER:O	2:I:87:ARG:NH1	2.36	0.57
1:B:331:ASN:HA	1:B:580:GLN:HA	1.86	0.57
2:H:91:THR:HG23	2:H:121:THR:HA	1.86	0.57
2:G:70:MET:SD	2:G:81:MET:HG2	2.44	0.57
4:D:365:THR:HG22	4:D:367:ASP:H	1.69	0.57
1:A:353:TRP:HZ3	1:A:355:ARG:HD3	1.69	0.57
1:A:102:ARG:HH21	1:A:122:ASN:HA	1.69	0.57
1:A:697:MET:HE3	1:B:864:LEU:HD11	1.86	0.57
1:C:776:LYS:NZ	1:C:780:GLU:OE2	2.38	0.57
1:A:328:ARG:HD3	1:A:531:THR:O	2.05	0.57
1:C:490:PHE:O	1:C:493:GLN:NE2	2.37	0.57
1:C:773:GLU:OE2	1:C:1019:ARG:NH1	2.35	0.57
2:I:12:LYS:HB2	2:I:122:VAL:HG22	1.87	0.57
1:C:189:LEU:HB3	1:C:208:THR:HB	1.87	0.57
3:K:91:GLN:NE2	3:K:100:SER:OG	2.37	0.57
1:B:353:TRP:HZ3	1:B:355:ARG:HD3	1.69	0.57
4:F:564:GLU:HB3	4:F:568:LEU:HD23	1.86	0.57
1:C:101:ILE:HG13	1:C:242:LEU:HD13	1.87	0.57
2:H:156:TYR:CE1	2:H:188:SER:HA	2.40	0.56
2:G:108:THR:OG1	3:J:36:HIS:NE2	2.34	0.56
4:F:48:TRP:CZ3	4:F:359:LEU:HB2	2.40	0.56
4:F:392:LEU:HD13	4:F:563:SER:HB3	1.87	0.56
4:F:245:ARG:NH2	4:F:603:PHE:O	2.38	0.56
4:D:116:LEU:HD11	4:D:187:LYS:HE2	1.86	0.56
1:B:577:ARG:HD3	1:B:582:LEU:HD22	1.86	0.56
1:B:645:THR:HG23	1:B:647:ALA:H	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:312:GLU:HG3	4:E:323:MET:HE1	1.86	0.56
1:A:34:ARG:NH1	1:A:191:GLU:OE2	2.38	0.56
4:F:394:ASN:OD1	4:F:401:HIS:NE2	2.38	0.56
1:A:34:ARG:NH2	1:A:219:GLY:O	2.38	0.56
1:B:905:ARG:NH1	1:B:1049:LEU:O	2.39	0.56
1:B:699:LEU:HB3	1:C:873:TYR:HE1	1.71	0.56
4:D:526:GLN:NE2	4:D:530:CYS:SG	2.78	0.56
2:I:5:VAL:HB	2:I:23:LYS:HB3	1.88	0.56
4:E:55:THR:OG1	4:E:58:ASN:ND2	2.37	0.55
1:A:909:ILE:HD12	1:A:1047:TYR:HB3	1.86	0.55
1:B:819:GLU:HG2	1:B:1054:GLN:HB3	1.87	0.55
2:H:69:MET:HB3	2:H:82:GLU:HB3	1.88	0.55
3:K:54:ASN:O	3:K:56:ARG:NH1	2.39	0.55
1:B:689:SER:OG	1:B:690:GLN:N	2.40	0.55
4:F:325:GLN:HA	4:F:328:TRP:HD1	1.71	0.55
2:I:86:LEU:HD12	2:I:122:VAL:HG21	1.88	0.55
1:B:566:GLY:N	1:B:575:ALA:O	2.39	0.55
2:H:39:GLN:OE1	3:L:40:GLN:NE2	2.40	0.55
3:K:152:ALA:HB3	3:K:199:GLN:HB2	1.88	0.55
1:B:258:TRP:HZ3	1:B:260:ALA:HB2	1.71	0.55
1:B:529:LYS:HE3	1:B:529:LYS:H	1.71	0.55
4:F:526:GLN:NE2	4:F:530:CYS:SG	2.79	0.55
2:I:91:THR:HG23	2:I:121:THR:HA	1.89	0.55
4:D:361:CYS:SG	4:D:363:LYS:NZ	2.80	0.55
1:A:328:ARG:NH2	1:A:533:LEU:HD13	2.22	0.55
1:B:329:PHE:HB3	1:B:330:PRO:HD3	1.89	0.55
4:F:408:MET:HA	4:F:411:SER:HB2	1.88	0.55
1:B:1076:THR:O	1:B:1097:SER:OG	2.24	0.54
1:C:172:SER:OG	1:C:173:GLN:N	2.40	0.54
1:C:1076:THR:O	1:C:1097:SER:OG	2.25	0.54
2:G:38:ARG:NH2	2:G:46:GLU:OE1	2.40	0.54
1:B:659:SER:OG	1:B:698:SER:OG	2.21	0.54
4:F:248:LEU:HB3	4:F:256:ILE:HD13	1.89	0.54
1:A:520:ALA:HB1	1:A:521:PRO:HD2	1.90	0.54
1:A:172:SER:OG	1:A:173:GLN:N	2.37	0.54
4:D:582:ARG:O	4:D:586:ASN:ND2	2.40	0.54
1:C:159:VAL:HG23	1:C:160:TYR:HD1	1.71	0.54
4:F:447:VAL:HA	4:F:450:LEU:HD12	1.88	0.54
1:C:983:ARG:HG3	1:C:984:LEU:HG	1.88	0.54
3:J:91:GLN:OE1	3:J:100:SER:OG	2.26	0.54
1:B:199:GLY:HA3	1:B:232:GLY:HA2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:VAL:HB	1:B:172:SER:HB3	1.90	0.54
1:B:490:PHE:O	1:B:493:GLN:NE2	2.41	0.54
4:E:83:TYR:O	4:E:101:GLN:NE2	2.41	0.54
3:J:6:GLN:NE2	3:J:90:CYS:SG	2.81	0.54
2:I:23:LYS:NZ	2:I:24:ALA:O	2.41	0.54
1:B:880:GLY:O	1:B:884:SER:OG	2.27	0.53
4:F:524:GLN:HB3	4:F:574:VAL:HG11	1.90	0.53
1:B:393:THR:HA	1:B:522:ALA:HA	1.91	0.53
1:B:731:MET:HB2	1:B:955:ASN:HD21	1.73	0.53
1:C:763:LEU:HD22	1:C:1008:VAL:HG21	1.91	0.53
1:A:97:LYS:HB2	1:A:186:PHE:HA	1.91	0.53
1:A:329:PHE:H	1:A:528:LYS:HZ3	1.55	0.53
4:E:392:LEU:HD13	4:E:563:SER:HB3	1.89	0.53
4:E:351:LEU:HB2	4:E:355:ASP:HB3	1.91	0.53
1:B:102:ARG:HG3	1:B:141:LEU:HD12	1.91	0.53
1:B:437:ASN:ND2	1:B:506:GLN:OE1	2.42	0.53
4:E:503:LEU:HD12	4:E:504:PHE:H	1.74	0.53
1:B:398:ASP:HB2	1:B:512:VAL:HB	1.91	0.53
4:F:83:TYR:O	4:F:101:GLN:NE2	2.41	0.53
4:F:233:ILE:HD13	4:F:450:LEU:HD13	1.90	0.53
4:F:482:ARG:NH2	4:F:613:TYR:OH	2.40	0.53
4:D:188:ASN:HD21	4:D:464:PHE:HA	1.74	0.53
2:G:2:VAL:HA	2:G:26:GLY:HA3	1.90	0.53
1:A:409:GLN:HB3	1:A:419:ALA:HB2	1.92	0.52
1:B:543:PHE:CD1	1:B:579:PRO:HD3	2.42	0.52
3:J:167:THR:HG1	3:J:180:SER:H	1.57	0.52
1:C:882:ILE:HG13	1:C:883:THR:HG23	1.91	0.52
2:G:28:SER:OG	2:G:98:ARG:NH1	2.42	0.52
1:A:494:SER:HB2	2:H:55:LEU:HD13	1.90	0.52
1:A:581:THR:HG23	1:A:583:GLU:HB2	1.91	0.52
1:C:880:GLY:O	1:C:884:SER:OG	2.26	0.52
3:L:63:ARG:NH1	3:L:79:GLY:O	2.42	0.52
4:D:524:GLN:HB3	4:D:574:VAL:HG11	1.92	0.52
1:B:134:GLN:HB2	1:B:161:SER:HB2	1.91	0.52
1:B:691:SER:OG	1:B:692:ILE:N	2.42	0.52
1:C:98:SER:HB3	1:C:181:GLY:HA2	1.91	0.52
1:C:102:ARG:HH21	1:C:122:ASN:HA	1.74	0.52
3:J:63:ARG:NH1	3:J:79:GLY:O	2.43	0.52
1:B:578:ASP:O	1:B:582:LEU:HD23	2.09	0.52
2:H:38:ARG:NH1	2:H:90:ASP:OD1	2.43	0.52
4:F:177:ARG:NH2	4:F:470:LYS:O	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:12:LYS:HE2	2:H:122:VAL:HG22	1.91	0.52
2:H:157:PHE:HB3	2:H:158:PRO:HD3	1.92	0.52
4:E:524:GLN:HB3	4:E:574:VAL:HG11	1.92	0.52
2:I:180:VAL:HG21	3:K:165:GLU:HB3	1.92	0.52
3:L:17:ARG:NH1	3:L:19:THR:OG1	2.43	0.52
4:E:144:LEU:HD12	4:E:148:LEU:HD22	1.91	0.52
4:E:213:ASP:OD1	4:E:213:ASP:N	2.43	0.52
1:A:142:ASP:HB2	1:A:156:GLU:HB2	1.92	0.52
1:C:457:ARG:NH1	1:C:459:SER:OG	2.43	0.52
4:F:144:LEU:HD12	4:F:148:LEU:HD22	1.92	0.52
1:A:327:VAL:HG22	1:A:542:ASN:HB3	1.91	0.52
1:A:359:SER:HA	1:A:524:VAL:HG23	1.91	0.52
1:A:393:THR:HG21	1:A:518:LEU:HB2	1.92	0.52
1:B:346:ARG:NH2	3:J:99:ASP:OD1	2.42	0.52
4:E:482:ARG:NH2	4:E:613:TYR:OH	2.41	0.52
1:B:872:GLN:O	1:B:875:SER:OG	2.26	0.51
1:C:327:VAL:HG22	1:C:542:ASN:HB3	1.90	0.51
1:A:327:VAL:HG12	1:A:528:LYS:NZ	2.25	0.51
1:A:804:GLN:NE2	1:A:935:GLN:OE1	2.43	0.51
1:A:363:ALA:HB2	1:A:524:VAL:HG13	1.92	0.51
1:B:487:ASN:ND2	4:E:24:GLN:OE1	2.42	0.51
1:C:355:ARG:NH2	1:C:464:PHE:CE1	2.77	0.51
1:C:981:LEU:HD21	1:C:993:ILE:HD11	1.92	0.51
4:E:50:TYR:HA	4:E:58:ASN:HB3	1.92	0.51
4:F:132:VAL:HB	4:F:144:LEU:HD21	1.91	0.51
1:A:102:ARG:HG3	1:A:141:LEU:HD12	1.93	0.51
1:C:346:ARG:NH2	3:K:99:ASP:OD1	2.44	0.51
1:A:236:THR:HG23	1:A:237:ARG:HG3	1.92	0.51
1:A:330:PRO:O	1:A:580:GLN:NE2	2.43	0.51
1:C:331:ASN:C	1:C:332:ILE:HD12	2.36	0.51
4:E:478:TRP:HE1	4:E:493:HIS:HB2	1.75	0.51
2:G:134:PRO:HA	2:G:151[A]:CYS:HA	1.93	0.51
2:G:134:PRO:HA	2:G:151[B]:CYS:HA	1.93	0.51
1:A:763:LEU:HD22	1:A:1008:VAL:HG21	1.92	0.51
4:E:552:GLN:NE2	4:E:556:ASN:OD1	2.44	0.51
1:C:134:GLN:HB2	1:C:161:SER:HB2	1.91	0.51
4:D:153:ALA:HA	4:D:277:ASN:HB3	1.93	0.51
1:B:37:TYR:OH	1:B:195:LYS:NZ	2.39	0.50
1:B:979:ASP:O	1:B:982:SER:OG	2.27	0.50
2:H:7:SER:HB3	2:H:21:SER:H	1.75	0.50
4:E:564:GLU:HB3	4:E:568:LEU:HD23	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:86:LEU:HB3	2:I:122:VAL:HG11	1.92	0.50
3:K:34:ASP:OD1	3:K:53:ASN:ND2	2.45	0.50
4:E:402:GLU:HB3	4:E:518:ARG:HD3	1.93	0.50
4:E:455:MET:HB2	4:E:480:MET:HE1	1.93	0.50
1:A:29:THR:HG23	1:A:62:VAL:HG13	1.92	0.50
1:B:701:ALA:HB3	1:C:787:GLN:HG2	1.93	0.50
4:E:32:PHE:HD1	4:E:76:GLN:HG3	1.76	0.50
2:G:10:GLU:HB2	2:G:120:VAL:HG22	1.93	0.50
1:C:966:LEU:HD22	1:C:1000:ARG:HH11	1.76	0.50
2:I:108:THR:OG1	3:K:36:HIS:NE2	2.38	0.50
1:A:328:ARG:HG2	1:A:530:SER:HA	1.93	0.50
1:B:159:VAL:HG23	1:B:160:TYR:HD1	1.74	0.50
1:B:528:LYS:HD2	1:B:528:LYS:N	2.27	0.50
4:F:597:ASP:O	4:F:600:LYS:NZ	2.41	0.50
1:C:599:THR:HG22	1:C:601:GLY:H	1.76	0.50
4:F:108:LEU:HD11	4:F:190:MET:HB3	1.93	0.50
4:E:261:CYS:HB2	4:E:488:VAL:HG23	1.93	0.50
1:B:98:SER:HB3	1:B:181:GLY:HA2	1.92	0.50
1:C:414:GLN:O	1:C:424:LYS:NZ	2.45	0.50
4:E:379:ILE:HG22	4:E:383:MET:HE1	1.94	0.50
2:G:91:THR:HG23	2:G:121:THR:HA	1.93	0.50
1:A:346:ARG:NH1	2:H:65:GLN:OE1	2.44	0.50
1:B:389:ASP:OD1	1:B:529:LYS:HD3	2.12	0.50
1:B:598:ILE:HB	1:B:609:ALA:HB3	1.94	0.50
1:C:365:TYR:HB2	1:C:388:ASN:HD21	1.76	0.50
2:H:2:VAL:HG21	2:H:98:ARG:HH12	1.77	0.50
4:F:171:GLU:O	4:F:175:GLN:NE2	2.37	0.50
4:F:552:GLN:NE2	4:F:556:ASN:OD1	2.45	0.50
1:A:594:GLY:H	1:A:613:GLN:CD	2.20	0.49
1:B:115:GLN:HE22	1:B:166:CYS:HA	1.76	0.49
1:C:383:SER:HB2	1:C:386:LYS:HB3	1.94	0.49
4:E:321:PRO:O	4:E:380:GLN:NE2	2.45	0.49
1:B:1000:ARG:O	1:B:1003:SER:OG	2.30	0.49
3:J:85:GLU:HG2	3:J:109:THR:HA	1.93	0.49
1:B:329:PHE:CD2	1:B:528:LYS:HG2	2.47	0.49
1:B:529:LYS:HE3	1:B:529:LYS:N	2.26	0.49
1:C:379:CYS:HA	1:C:432:CYS:HA	1.94	0.49
4:E:431:ASP:OD1	4:E:431:ASP:N	2.45	0.49
2:G:86:LEU:HD12	2:G:122:VAL:HG21	1.94	0.49
1:B:879:ALA:O	1:B:883:THR:OG1	2.25	0.49
1:B:974:SER:HB3	1:B:980:ILE:HG23	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:329:PHE:HE2	1:C:528:LYS:HD3	1.78	0.49
1:C:375:PHE:HE2	1:C:407:VAL:HG21	1.77	0.49
4:F:111:ASP:OD1	4:F:111:ASP:N	2.45	0.49
1:A:328:ARG:NH1	1:A:533:LEU:HB2	2.26	0.49
1:A:689:SER:OG	1:A:690:GLN:N	2.45	0.49
1:B:936:ASP:O	1:B:939:SER:OG	2.28	0.49
3:L:6:GLN:NE2	3:L:90:CYS:SG	2.86	0.49
2:I:149:LEU:HD12	2:I:222:VAL:HG22	1.94	0.49
4:D:252:TYR:HB3	4:D:255:TYR:HD2	1.76	0.49
1:B:319:ARG:HH21	1:C:740:MET:HB2	1.78	0.49
4:F:431:ASP:OD1	4:F:431:ASP:N	2.43	0.49
1:A:342:PHE:HE1	1:A:511:VAL:HG11	1.77	0.49
4:E:476:LYS:O	4:E:480:MET:HG3	2.12	0.49
1:A:253:ASP:OD1	1:A:253:ASP:N	2.45	0.49
3:J:141:ILE:HD13	3:J:149:VAL:HG11	1.95	0.49
3:K:20:ILE:HB	3:K:75:LEU:HB3	1.95	0.49
1:A:329:PHE:H	1:A:528:LYS:NZ	2.11	0.49
1:A:703:ASN:ND2	1:B:787:GLN:OE1	2.39	0.49
1:B:334:ASN:HB2	1:B:362:VAL:CB	2.43	0.49
1:C:102:ARG:HG3	1:C:141:LEU:HD12	1.94	0.49
3:L:48:LEU:HD21	3:L:51:TYR:HB3	1.95	0.49
4:E:526:GLN:NE2	4:E:530:CYS:SG	2.83	0.49
2:G:4:LEU:HD23	2:G:24:ALA:HA	1.94	0.49
1:A:246:ARG:HD3	1:A:252:GLY:HA2	1.94	0.49
2:G:149:LEU:HD12	2:G:222:VAL:HG22	1.95	0.49
3:K:14:PRO:HD3	3:K:110:VAL:HG13	1.95	0.49
1:A:195:LYS:HB2	1:A:202:LYS:HB2	1.95	0.48
1:A:966:LEU:O	1:A:975:SER:OG	2.30	0.48
4:E:346:PRO:HA	4:E:360:MET:HB3	1.95	0.48
2:G:70:MET:HE1	2:G:81:MET:HG2	1.96	0.48
2:I:98:ARG:HE	2:I:112:ASP:HB3	1.77	0.48
4:D:213:ASP:OD1	4:D:213:ASP:N	2.45	0.48
1:A:116:SER:N	1:A:131:CYS:O	2.46	0.48
1:B:355:ARG:O	3:J:95:SER:OG	2.28	0.48
2:H:67:ARG:HH11	2:H:83:LEU:HD11	1.78	0.48
1:A:365:TYR:HD1	1:A:368:LEU:HD11	1.77	0.48
1:A:697:MET:CE	1:B:864:LEU:HD11	2.43	0.48
1:C:175:PHE:HE2	1:C:190:ARG:HH12	1.61	0.48
4:F:261:CYS:HB2	4:F:488:VAL:HG23	1.94	0.48
3:K:20:ILE:HD12	3:K:75:LEU:HD23	1.95	0.48
1:B:246:ARG:HH22	1:B:254:SER:HB3	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:60:TYR:HE1	2:G:70:MET:HB2	1.77	0.48
1:A:355:ARG:O	3:L:95:SER:OG	2.26	0.48
1:C:93:ALA:HA	1:C:191:GLU:HA	1.95	0.48
4:E:55:THR:HG1	4:E:58:ASN:HD22	1.60	0.48
4:E:168:TRP:HD1	4:E:169:ARG:HH11	1.61	0.48
4:F:582:ARG:O	4:F:586:ASN:ND2	2.47	0.48
1:C:185:ASN:ND2	1:C:211:ASN:OD1	2.40	0.48
4:F:168:TRP:CD1	4:F:270:MET:HE1	2.49	0.48
4:F:365:THR:HG22	4:F:367:ASP:H	1.79	0.48
1:B:337:PRO:HD2	1:B:358:ILE:HG23	1.95	0.48
1:C:453:TYR:HE2	1:C:455:LEU:HD13	1.79	0.48
2:G:48:MET:HE2	2:G:81:MET:SD	2.54	0.48
3:J:171:LYS:NZ	3:J:175:ASN:O	2.47	0.48
4:F:360:MET:SD	4:F:360:MET:N	2.87	0.48
1:A:334:ASN:C	1:A:335:LEU:HG	2.39	0.48
1:C:656:VAL:HG22	1:C:658:ASN:H	1.78	0.48
2:H:128:LYS:NZ	2:H:155:ASP:HB3	2.29	0.48
2:G:38:ARG:NH1	2:G:90:ASP:OD1	2.47	0.48
4:F:109:SER:OG	4:F:111:ASP:OD1	2.29	0.48
3:K:154:LYS:HE2	3:K:157:SER:H	1.79	0.48
1:B:34:ARG:NH2	1:B:219:GLY:O	2.46	0.48
4:F:331:SER:OG	4:F:358:ILE:O	2.27	0.48
1:C:164:ASN:OD1	1:C:165:ASN:N	2.43	0.48
1:A:328:ARG:HH22	1:A:533:LEU:HD13	1.79	0.47
1:A:490:PHE:O	1:A:493:GLN:NE2	2.46	0.47
1:B:142:ASP:HB2	1:B:156:GLU:HB2	1.95	0.47
1:C:126:VAL:HB	1:C:172:SER:HB3	1.96	0.47
3:J:154:LYS:HD2	3:J:159:PRO:HB3	1.96	0.47
1:B:578:ASP:OD2	1:B:583:GLU:HB3	2.14	0.47
1:C:105:ILE:HG23	1:C:241:LEU:HD11	1.97	0.47
1:C:412:PRO:HB3	1:C:427:ASP:HA	1.96	0.47
4:E:588:PHE:O	4:E:592:PHE:N	2.45	0.47
3:J:137:LEU:HD12	3:J:183:LEU:HD23	1.95	0.47
1:A:360:ASN:N	1:A:523:THR:HB	2.19	0.47
1:C:747:THR:O	1:C:750:SER:OG	2.28	0.47
4:F:297:MET:HB2	4:F:302:TRP:HB2	1.95	0.47
4:D:55:THR:HG1	4:D:58:ASN:HD22	1.60	0.47
2:I:150:GLY:HA2	2:I:165:TRP:HH2	1.79	0.47
4:D:229:THR:HB	4:D:581:VAL:HG13	1.97	0.47
1:A:402:ILE:HB	1:A:406:GLU:HG3	1.95	0.47
1:A:465:GLU:HA	2:H:105:TYR:HA	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:737:ASP:HB2	1:B:740:MET:HB2	1.97	0.47
2:H:125:ALA:HB3	2:H:157:PHE:CE1	2.49	0.47
2:I:158:PRO:HD2	2:I:211:HIS:CE1	2.49	0.47
1:A:442:ASP:O	1:A:448:ASN:ND2	2.39	0.47
1:B:661:GLU:OE2	1:B:662:CYS:N	2.48	0.47
3:J:4:LEU:HB2	3:J:103:GLY:HA2	1.96	0.47
4:F:528:ALA:HB2	4:F:574:VAL:HG12	1.96	0.47
4:D:232:GLU:OE1	4:D:582:ARG:NH2	2.44	0.47
1:A:335:LEU:HD22	1:A:364:ASP:CB	2.39	0.47
1:B:346:ARG:NH1	2:G:65:GLN:OE1	2.48	0.47
1:B:557:LYS:O	1:B:584:ILE:HG21	2.14	0.47
1:C:974:SER:HB2	1:C:983:ARG:HH22	1.79	0.47
4:F:112:LYS:NZ	4:F:189:GLU:OE2	2.46	0.47
4:F:517:THR:HG22	4:F:579:MET:HE1	1.97	0.47
4:D:82:MET:SD	4:D:82:MET:N	2.83	0.47
1:A:531:THR:O	1:A:532:ASN:C	2.57	0.47
1:A:826:VAL:HG11	1:A:1057:PRO:HG2	1.96	0.47
1:A:152:TRP:HB2	1:A:179:LEU:HB3	1.96	0.47
1:C:189:LEU:N	1:C:208:THR:O	2.47	0.47
1:C:708:SER:OG	1:C:709:ASN:N	2.48	0.47
1:C:948:LEU:HD21	1:C:1059:GLY:HA3	1.97	0.47
4:E:582:ARG:O	4:E:586:ASN:ND2	2.48	0.47
2:I:48:MET:HE2	2:I:81:MET:HE1	1.96	0.47
1:C:936:ASP:O	1:C:939:SER:OG	2.29	0.46
4:D:205:GLY:HA2	4:D:219:ARG:HD2	1.96	0.46
1:A:916:LEU:HD12	1:A:923:ILE:HD12	1.97	0.46
1:A:1030:SER:HA	1:A:1034:LEU:HD12	1.97	0.46
1:B:117:LEU:HG	1:B:130:VAL:HG22	1.97	0.46
1:B:125:ASN:OD1	1:B:126:VAL:N	2.48	0.46
2:I:60:TYR:HE1	2:I:70:MET:HG2	1.80	0.46
1:A:117:LEU:HG	1:A:130:VAL:HG22	1.98	0.46
1:A:534:VAL:CG1	1:A:539:VAL:HG21	2.44	0.46
1:B:502:GLY:HA3	4:E:354:GLY:HA3	1.97	0.46
1:B:738:CYS:O	1:B:742:ILE:N	2.42	0.46
1:C:503:VAL:HA	1:C:506:GLN:HG3	1.97	0.46
4:D:389:PRO:HD2	4:D:392:LEU:HD12	1.96	0.46
1:B:578:ASP:O	1:B:582:LEU:HA	2.15	0.46
4:E:132:VAL:HB	4:E:144:LEU:HD21	1.96	0.46
1:A:396:TYR:HB2	1:A:514:SER:HB3	1.98	0.46
1:B:1083:HIS:ND1	1:B:1084:ASP:OD1	2.48	0.46
1:C:715:PRO:HA	1:C:1072:GLU:HA	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:237:TYR:OH	4:F:485:VAL:O	2.34	0.46
2:I:72:ALA:HA	2:I:79:ALA:HA	1.98	0.46
3:K:149:VAL:HG13	3:K:200:VAL:HG13	1.97	0.46
4:D:457:GLU:HG2	4:D:512:PHE:HB2	1.97	0.46
1:A:335:LEU:HA	1:A:362:VAL:O	2.16	0.46
4:E:365:THR:HG22	4:E:367:ASP:H	1.80	0.46
4:E:505:HIS:HB3	4:E:510:TYR:HB2	1.98	0.46
1:A:199:GLY:HA3	1:A:232:GLY:HA2	1.96	0.46
1:B:153:MET:HE1	1:B:155:SER:HB3	1.97	0.46
1:B:543:PHE:HD1	1:B:579:PRO:CD	2.28	0.46
1:C:81:ASN:HB3	1:C:239:GLN:HE21	1.80	0.46
1:A:29:THR:HB	1:A:214:ARG:HH22	1.80	0.46
1:A:725:GLU:OE1	1:A:1064:HIS:NE2	2.45	0.46
1:B:280:ASN:OD1	1:B:284:THR:N	2.45	0.46
2:G:12:LYS:HB2	2:G:122:VAL:HG22	1.97	0.46
2:G:23:LYS:NZ	2:G:24:ALA:O	2.45	0.46
4:F:80:ALA:O	4:F:101:GLN:NE2	2.37	0.46
4:F:152:MET:HG3	4:F:270:MET:HB3	1.98	0.46
4:D:503:LEU:HD12	4:D:504:PHE:H	1.80	0.46
1:B:856:ASN:HD21	1:B:963:VAL:HG12	1.81	0.46
1:C:355:ARG:NH2	1:C:464:PHE:HE1	2.13	0.46
1:C:437:ASN:ND2	1:C:506:GLN:OE1	2.48	0.46
2:H:127:THR:HG1	2:H:157:PHE:HD2	1.63	0.46
4:E:245:ARG:NH2	4:E:605:GLY:O	2.47	0.46
4:F:55:THR:OG1	4:F:58:ASN:OD1	2.33	0.46
4:D:230:PHE:HA	4:D:233:ILE:HD12	1.97	0.46
1:A:95:THR:HA	1:A:189:LEU:HA	1.98	0.46
1:A:153:MET:HE1	1:A:155:SER:HB3	1.98	0.46
1:A:776:LYS:NZ	1:A:780:GLU:OE2	2.48	0.46
1:A:787:GLN:OE1	1:C:703:ASN:ND2	2.41	0.46
1:B:32:PHE:HB3	1:B:218:GLN:HG3	1.97	0.46
4:E:447:VAL:HA	4:E:450:LEU:HD12	1.98	0.46
3:J:125:PRO:HB3	3:J:129:GLU:HG2	1.97	0.46
4:F:582:ARG:HE	4:F:585:LEU:HD12	1.81	0.46
2:I:102:TYR:OH	3:K:34:ASP:OD2	2.34	0.46
1:C:342:PHE:HE2	1:C:368:LEU:HD22	1.81	0.45
3:L:186:THR:HG22	3:L:188:GLU:H	1.81	0.45
4:E:560:LEU:O	4:E:563:SER:OG	2.32	0.45
4:F:398:GLU:OE1	4:F:514:ARG:NH2	2.49	0.45
4:D:79:LEU:HA	4:D:82:MET:HE1	1.98	0.45
4:D:574:VAL:HG23	4:D:576:ALA:H	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:879:ALA:O	1:C:883:THR:OG1	2.31	0.45
4:D:29:LEU:O	4:D:33:ASN:ND2	2.49	0.45
1:A:792:PRO:HG3	1:C:707:TYR:HB3	1.99	0.45
2:G:39:GLN:OE1	3:J:40:GLN:NE2	2.47	0.45
2:I:132:VAL:HG22	2:I:153:VAL:HG22	1.98	0.45
4:D:586:ASN:HA	4:D:589:GLU:HB2	1.98	0.45
1:A:29:THR:OG1	1:A:30:ASN:N	2.49	0.45
1:A:469:SER:HA	2:H:101:GLY:HA3	1.98	0.45
1:B:773:GLU:OE2	1:B:1019:ARG:NH1	2.41	0.45
1:C:528:LYS:HD2	1:C:528:LYS:H	1.81	0.45
3:J:41:LEU:HB2	3:J:44:ALA:HB3	1.98	0.45
4:F:398:GLU:HG3	4:F:514:ARG:HB3	1.98	0.45
4:D:315:PHE:HD1	4:D:320:LEU:HD12	1.81	0.45
1:A:461:LEU:HD21	1:A:467:ASP:HB2	1.98	0.45
1:A:721:SER:HG	1:A:1066:THR:HG1	1.64	0.45
1:B:48:LEU:HB3	1:B:276:LEU:HD11	1.99	0.45
1:B:439:ASN:O	1:B:443:SER:OG	2.30	0.45
1:C:919:ASN:HB2	1:C:922:LEU:HB3	1.99	0.45
2:H:39:GLN:HB2	2:H:45:LEU:HD23	1.98	0.45
2:G:179:ALA:HB1	2:G:187:TYR:HB3	1.97	0.45
3:K:145:TYR:HA	3:K:177:TYR:HD1	1.82	0.45
2:H:111:MET:SD	2:H:111:MET:N	2.90	0.45
2:H:174:VAL:HA	2:H:193:VAL:HG12	1.99	0.45
2:H:211:HIS:CD2	2:H:213:PRO:HD2	2.52	0.45
4:F:460:ARG:NH2	4:F:510:TYR:O	2.45	0.45
4:D:303:ASP:N	4:D:303:ASP:OD1	2.50	0.45
1:B:91:TYR:N	1:B:268:GLY:O	2.48	0.45
1:B:645:THR:OG1	1:B:646:ARG:N	2.50	0.45
1:C:342:PHE:HE1	1:C:511:VAL:HG11	1.82	0.45
1:C:373:PRO:HB2	1:C:436:TRP:HB2	1.97	0.45
1:C:396:TYR:HB2	1:C:514:SER:HB3	1.98	0.45
1:C:528:LYS:HD2	1:C:528:LYS:N	2.32	0.45
4:E:177:ARG:NH2	4:E:470:LYS:O	2.50	0.45
4:E:226:VAL:HG21	4:E:513:ILE:HD11	1.99	0.45
4:F:168:TRP:HD1	4:F:169:ARG:HH11	1.65	0.45
3:K:169:PRO:HA	3:K:179:ALA:HB2	1.98	0.45
3:L:65:SER:N	3:L:76:ALA:O	2.48	0.45
4:E:166:GLU:OE2	4:E:170:SER:OG	2.34	0.45
2:G:39:GLN:HB2	2:G:45:LEU:HD23	1.97	0.45
4:D:482:ARG:O	4:D:606:TRP:NE1	2.50	0.45
1:C:34:ARG:NH2	1:C:219:GLY:O	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:ILE:HD12	1:C:110:LEU:HD13	1.99	0.45
1:C:258:TRP:HZ3	1:C:260:ALA:HB2	1.82	0.45
1:C:742:ILE:HA	1:C:1000:ARG:HD2	1.98	0.45
1:C:902:MET:HE1	1:C:1049:LEU:HD13	1.99	0.45
2:I:36:TRP:HB2	2:I:48:MET:HE3	1.98	0.45
3:K:194:ARG:HH12	3:K:212:ALA:HB1	1.82	0.45
1:B:721:SER:OG	1:B:1066:THR:OG1	2.36	0.44
1:C:439:ASN:O	1:C:443:SER:OG	2.31	0.44
4:E:26:LYS:HE2	4:E:90:ASN:HD22	1.81	0.44
4:F:378:HIS:CE1	4:F:401:HIS:HB3	2.52	0.44
1:B:435:ALA:HB2	1:B:510:VAL:HG13	1.98	0.44
2:H:181:LEU:HD13	2:H:187:TYR:HE1	1.83	0.44
4:E:269:ASP:OD1	4:E:269:ASP:N	2.50	0.44
4:E:586:ASN:HA	4:E:589:GLU:HB2	1.99	0.44
1:A:617:CYS:N	1:A:649:CYS:SG	2.90	0.44
1:B:356:LYS:HB3	1:B:397:ALA:HB3	2.00	0.44
1:B:383:SER:HB2	1:B:386:LYS:HB3	1.99	0.44
1:C:318:PHE:HD2	1:C:615:VAL:HG21	1.81	0.44
2:G:85:SER:HB3	2:G:87:ARG:NH1	2.32	0.44
3:K:107:LYS:NZ	3:K:108:VAL:O	2.47	0.44
4:D:166:GLU:OE2	4:D:170:SER:OG	2.36	0.44
1:B:110:LEU:HB3	1:B:135:PHE:HD2	1.82	0.44
1:C:887:THR:HG21	1:C:894:LEU:HD12	1.99	0.44
2:H:12:LYS:HE3	2:H:86:LEU:HG	2.00	0.44
3:J:40:GLN:HB3	3:J:87:ASP:HB2	1.98	0.44
4:D:529:LEU:HD21	4:D:554:LEU:HB2	1.99	0.44
1:A:331:ASN:C	1:A:332:ILE:HD12	2.43	0.44
1:B:662:CYS:SG	1:B:697:MET:HG3	2.56	0.44
3:J:121:THR:HB	3:J:140:LEU:HB2	1.99	0.44
4:F:503:LEU:HB3	4:F:506:VAL:HG22	2.00	0.44
1:A:37:TYR:HB3	1:A:223:LEU:HD12	1.99	0.44
1:A:328:ARG:HG3	1:A:531:THR:OG1	2.17	0.44
1:A:365:TYR:HB2	1:A:388:ASN:HD21	1.83	0.44
1:C:138:ASP:N	1:C:138:ASP:OD1	2.49	0.44
4:E:457:GLU:OE2	4:E:460:ARG:NH2	2.50	0.44
3:K:40:GLN:HA	3:K:46:PRO:HA	2.00	0.44
4:D:488:VAL:HG11	4:D:611:SER:HA	2.00	0.44
1:A:970:PHE:HA	1:B:756:TYR:HE1	1.82	0.44
1:A:1076:THR:O	1:A:1097:SER:OG	2.35	0.44
1:B:398:ASP:OD2	1:B:423:TYR:OH	2.32	0.44
2:G:18:VAL:HG12	2:G:86:LEU:HD21	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:VAL:HG21	1:A:418:ILE:HG23	1.99	0.44
1:C:425:LEU:HD21	1:C:512:VAL:HG11	2.00	0.44
4:E:232:GLU:OE1	4:E:582:ARG:NH2	2.48	0.44
4:E:480:MET:HE2	4:E:484:ILE:HD12	1.99	0.44
2:G:98:ARG:HE	2:G:112:ASP:HB3	1.83	0.44
2:G:127:THR:HA	2:G:157:PHE:HB3	2.00	0.44
2:G:206:ILE:HG22	2:G:221:LYS:HA	2.00	0.44
4:F:325:GLN:O	4:F:328:TRP:N	2.51	0.44
1:C:529:LYS:HD2	1:C:529:LYS:N	2.33	0.44
4:E:378:HIS:HE1	4:E:401:HIS:HB3	1.83	0.44
4:E:582:ARG:HE	4:E:585:LEU:HD12	1.82	0.44
3:K:199:GLN:HE22	3:K:208:GLU:HG2	1.82	0.44
4:D:390:PHE:HA	4:D:393:ARG:HD3	2.00	0.44
1:C:353:TRP:CD1	1:C:353:TRP:H	2.35	0.43
1:C:355:ARG:NH2	1:C:398:ASP:OD1	2.45	0.43
1:C:128:ILE:HD13	1:C:170:TYR:HD2	1.84	0.43
4:F:208:GLU:OE2	4:F:210:ASN:ND2	2.51	0.43
4:D:402:GLU:HB3	4:D:518:ARG:HD3	1.98	0.43
1:A:442:ASP:OD1	1:A:509:ARG:NH2	2.51	0.43
1:B:328:ARG:HD3	1:B:531:THR:OG1	2.18	0.43
1:B:365:TYR:HD1	1:B:368:LEU:HD11	1.81	0.43
1:B:490:PHE:CE2	1:B:492:LEU:HB2	2.54	0.43
2:H:41:PRO:HD3	2:H:92:ALA:HA	2.00	0.43
3:L:37:TRP:HB2	3:L:50:ILE:HB	2.00	0.43
3:K:82:VAL:HA	3:K:110:VAL:HG21	2.01	0.43
1:A:326:ILE:HB	1:A:531:THR:OG1	2.17	0.43
1:A:359:SER:HB3	1:A:523:THR:CB	2.48	0.43
1:B:776:LYS:NZ	1:B:780:GLU:OE2	2.51	0.43
1:C:355:ARG:HD3	1:C:356:LYS:N	2.34	0.43
2:H:179:ALA:HB1	2:H:187:TYR:HB3	1.99	0.43
3:J:90:CYS:O	3:J:103:GLY:N	2.52	0.43
1:B:418:ILE:HA	1:B:422:ASN:HB2	2.00	0.43
4:E:271:TRP:CD2	4:E:503:LEU:HD13	2.54	0.43
2:G:36:TRP:CD1	2:G:81:MET:HG3	2.53	0.43
3:K:48:LEU:HD11	3:K:51:TYR:HB3	1.99	0.43
1:A:67:ALA:HB3	1:A:263:ALA:HB3	2.00	0.43
1:A:395:VAL:HG21	1:A:524:VAL:HG11	2.01	0.43
1:A:1092:GLU:O	1:A:1107:ARG:NH2	2.52	0.43
1:C:360:ASN:H	1:C:523:THR:HB	1.84	0.43
4:F:315:PHE:HA	4:F:318:VAL:HG22	2.01	0.43
4:F:341:LYS:HA	4:F:341:LYS:HD3	1.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:389:PRO:HD2	4:F:392:LEU:HD12	1.99	0.43
1:A:895:GLN:NE2	1:C:1074:ASN:OD1	2.52	0.43
3:L:94:ASP:HB3	3:L:97:LEU:HB2	2.01	0.43
4:D:198:ASP:N	4:D:198:ASP:OD1	2.49	0.43
1:B:365:TYR:HB2	1:B:388:ASN:HD21	1.83	0.43
4:E:535:HIS:NE2	4:E:538:PRO:O	2.49	0.43
4:E:597:ASP:O	4:E:600:LYS:NZ	2.38	0.43
4:F:462:MET:HB3	4:F:468:ILE:HG13	2.00	0.43
4:F:523:PHE:HB3	4:F:583:PRO:HB2	2.00	0.43
2:I:178:PRO:HG3	3:K:170:SER:HB3	2.01	0.43
1:C:282:ASN:CG	6:C:1301:NAG:HN2	2.25	0.43
2:H:34:ILE:HB	2:H:51:ILE:HG23	2.01	0.43
4:F:201:ASP:OD2	4:F:219:ARG:NH1	2.52	0.43
1:C:616:ASN:HA	1:C:644:GLN:HE22	1.84	0.42
1:C:886:TRP:HZ3	1:C:905:ARG:HD3	1.84	0.42
4:E:455:MET:HG2	4:E:485:VAL:HG23	2.01	0.42
4:F:288:LYS:HD2	4:F:433:GLU:HB2	2.00	0.42
1:B:354:ASN:ND2	3:J:95:SER:O	2.51	0.42
4:E:528:ALA:HB2	4:E:574:VAL:HG12	2.00	0.42
2:G:163:VAL:HG22	2:G:209:VAL:HG12	2.00	0.42
4:D:394:ASN:OD1	4:D:401:HIS:NE2	2.52	0.42
1:B:568:ASP:OD1	1:B:572:THR:N	2.47	0.42
1:C:375:PHE:N	1:C:435:ALA:O	2.52	0.42
4:F:32:PHE:HD1	4:F:76:GLN:HG3	1.82	0.42
4:D:267:LEU:HD11	4:D:487:VAL:HG21	2.00	0.42
1:A:418:ILE:HA	1:A:422:ASN:HB2	2.00	0.42
1:B:764:LYS:O	1:B:768:THR:OG1	2.32	0.42
4:F:92:THR:HB	4:F:392:LEU:HD11	2.02	0.42
2:I:135:LEU:N	2:I:150:GLY:O	2.46	0.42
4:D:109:SER:OG	4:D:111:ASP:OD1	2.31	0.42
1:A:102:ARG:NH2	1:A:122:ASN:OD1	2.53	0.42
1:B:214:ARG:O	1:B:215:ASP:HB3	2.20	0.42
4:E:169:ARG:HG3	4:E:499:ASP:OD1	2.19	0.42
2:G:134:PRO:HD3	2:G:220:LYS:HG2	2.00	0.42
4:F:166:GLU:OE2	4:F:170:SER:OG	2.37	0.42
4:F:312:GLU:HG3	4:F:323:MET:HE3	2.01	0.42
4:D:73:LEU:HD23	4:D:73:LEU:HA	1.91	0.42
1:A:328:ARG:HG3	1:A:531:THR:HG1	1.85	0.42
1:A:482:GLY:O	2:H:31:ASN:ND2	2.52	0.42
1:B:856:ASN:HD22	1:B:966:LEU:HD11	1.85	0.42
1:C:86:PHE:H	1:C:237:ARG:HA	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:106:PHE:HB2	1:C:117:LEU:HB2	2.01	0.42
3:K:137:LEU:HD12	3:K:183:LEU:HD23	2.02	0.42
1:A:66:HIS:HB2	1:A:78:ARG:HG2	2.02	0.42
1:A:598:ILE:HB	1:A:609:ALA:HB3	2.01	0.42
1:A:877:LEU:HD23	1:A:877:LEU:HA	1.89	0.42
3:L:65:SER:HB3	3:L:76:ALA:HB3	2.00	0.42
4:E:80:ALA:O	4:E:101:GLN:NE2	2.39	0.42
4:E:233:ILE:HD13	4:E:450:LEU:HD13	2.01	0.42
2:G:78:THR:HG23	2:G:80:TYR:CE2	2.54	0.42
4:F:116:LEU:HD11	4:F:187:LYS:HE2	2.02	0.42
4:F:248:LEU:HD12	4:F:262:LEU:HD22	2.02	0.42
1:B:791:THR:HG21	1:B:806:LEU:HD22	2.01	0.42
1:C:97:LYS:HD3	1:C:187:LYS:H	1.85	0.42
2:H:154:LYS:HD2	2:H:188:SER:CB	2.47	0.42
4:E:488:VAL:HG11	4:E:611:SER:HA	2.01	0.42
1:A:125:ASN:HA	1:A:174:PRO:HA	2.02	0.42
1:C:448:ASN:HB3	1:C:497:PHE:HB2	2.01	0.42
4:E:323:MET:N	4:E:323:MET:SD	2.93	0.42
4:D:482:ARG:HD2	4:D:488:VAL:HG22	2.00	0.42
1:B:713:ALA:HB2	1:C:895:GLN:HG2	2.02	0.42
1:A:1018:ILE:HD13	1:A:1018:ILE:HA	1.92	0.41
1:C:398:ASP:HB2	1:C:512:VAL:HB	2.01	0.41
4:E:380:GLN:HA	4:E:383:MET:HE2	2.01	0.41
4:F:248:LEU:HD23	4:F:248:LEU:HA	1.92	0.41
3:K:94:ASP:HB3	3:K:97:LEU:HB2	2.02	0.41
1:B:66:HIS:HB2	1:B:78:ARG:HG2	2.02	0.41
1:B:290:ASP:OD1	1:B:291:CYS:N	2.53	0.41
1:C:355:ARG:O	3:K:95:SER:OG	2.36	0.41
1:C:355:ARG:CD	1:C:396:TYR:HD1	2.23	0.41
4:F:97:LEU:HD23	4:F:97:LEU:HA	1.92	0.41
1:B:342:PHE:HE1	1:B:511:VAL:HG11	1.86	0.41
4:E:474:MET:SD	4:E:478:TRP:NE1	2.92	0.41
2:G:150:GLY:HA2	2:G:165:TRP:HH2	1.85	0.41
1:A:534:VAL:HG13	1:A:539:VAL:HG21	2.03	0.41
1:C:40:ASP:N	1:C:40:ASP:OD1	2.53	0.41
4:F:90:ASN:CG	4:F:93:VAL:HG23	2.45	0.41
4:F:161:ARG:HB2	4:F:265:HIS:HD2	1.85	0.41
4:F:492:PRO:HD3	4:F:613:TYR:CG	2.55	0.41
1:A:670:ILE:N	1:A:697:MET:HE2	2.35	0.41
1:C:328:ARG:HH12	1:C:533:LEU:HB2	1.85	0.41
2:H:129:GLY:HA2	2:H:130:PRO:HD3	1.92	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:211:HIS:CE1	2:H:214:SER:HG	2.31	0.41
4:E:457:GLU:OE2	4:E:512:PHE:N	2.40	0.41
3:K:128:GLU:HB2	3:K:131:GLN:HE21	1.84	0.41
4:D:505:HIS:HB3	4:D:510:TYR:HB2	2.02	0.41
1:A:359:SER:HB3	1:A:523:THR:HB	2.03	0.41
1:A:747:THR:O	1:A:750:SER:OG	2.31	0.41
1:C:355:ARG:O	3:K:31:ALA:HB1	2.21	0.41
2:H:110:TYR:HB3	3:L:36:HIS:CE1	2.55	0.41
2:G:70:MET:CE	2:G:81:MET:HG2	2.51	0.41
2:I:163:VAL:HG22	2:I:209:VAL:HG12	2.03	0.41
1:A:333:THR:HB	1:A:335:LEU:HD23	2.02	0.41
1:A:428:ASP:OD1	1:A:428:ASP:N	2.50	0.41
4:F:535:HIS:NE2	4:F:538:PRO:O	2.49	0.41
4:D:171:GLU:O	4:D:175:GLN:NE2	2.53	0.41
1:B:543:PHE:C	1:B:544:ASN:CG	2.89	0.41
1:C:44:ARG:HD2	1:C:279:TYR:CZ	2.55	0.41
3:J:172:GLN:HE21	3:J:176:LYS:HB2	1.85	0.41
1:A:294:ASP:O	1:A:298:GLU:N	2.49	0.41
1:A:568:ASP:OD1	1:A:572:THR:N	2.45	0.41
1:B:379:CYS:HA	1:B:432:CYS:HA	2.02	0.41
1:B:878:LEU:HD13	1:B:878:LEU:HA	1.91	0.41
1:B:966:LEU:O	1:B:975:SER:OG	2.38	0.41
1:B:1046:GLY:HA2	1:C:890:ALA:HB1	2.02	0.41
3:L:141:ILE:HD13	3:L:149:VAL:HG11	2.02	0.41
1:A:710:ASN:OD1	1:A:710:ASN:N	2.54	0.41
4:E:472:GLN:NE2	4:E:476:LYS:HB2	2.36	0.41
4:F:47:SER:HB3	4:F:349:TRP:HZ3	1.86	0.41
4:F:488:VAL:HG11	4:F:611:SER:HA	2.02	0.41
2:I:146:THR:HA	2:I:196:PRO:HA	2.03	0.41
1:A:117:LEU:HD11	1:A:233:ILE:HD11	2.03	0.40
1:A:336:CYS:N	1:A:362:VAL:O	2.54	0.40
4:E:144:LEU:HG	4:E:168:TRP:CE3	2.56	0.40
4:F:144:LEU:HG	4:F:168:TRP:CE3	2.56	0.40
2:I:68:VAL:HG22	2:I:83:LEU:HD13	2.02	0.40
2:I:150:GLY:HA2	2:I:165:TRP:CH2	2.55	0.40
3:K:165:GLU:O	3:K:182:TYR:N	2.40	0.40
4:D:453:THR:HG23	4:D:512:PHE:HB3	2.03	0.40
4:D:455:MET:O	4:D:459:TRP:N	2.48	0.40
1:B:328:ARG:HH12	1:B:533:LEU:HB2	1.85	0.40
1:C:355:ARG:HH22	1:C:464:PHE:HD1	1.60	0.40
1:C:440:LYS:HA	1:C:440:LYS:HD3	1.82	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:94:TYR:O	2:G:118:THR:N	2.49	0.40
3:J:27:SER:HB2	3:J:95:SER:H	1.87	0.40
4:F:47:SER:HA	4:F:62:MET:HE1	2.04	0.40
4:F:105:SER:O	4:F:113:SER:OG	2.40	0.40
2:I:150:GLY:HA3	2:I:192:VAL:HG22	2.03	0.40
1:A:353:TRP:CZ3	1:A:355:ARG:HD3	2.53	0.40
1:B:317:ASN:HA	1:B:594:GLY:HA2	2.04	0.40
1:C:611:LEU:HD22	1:C:666:ILE:HG23	2.03	0.40
4:E:583:PRO:O	4:E:587:TYR:HB2	2.22	0.40
3:K:90:CYS:O	3:K:103:GLY:N	2.55	0.40
4:D:158:TYR:HE1	4:D:266:LEU:HD11	1.86	0.40
4:D:341:LYS:HA	4:D:341:LYS:HD3	1.88	0.40
1:A:328:ARG:NH1	1:A:580:GLN:HG3	2.37	0.40
1:B:210:ILE:HD12	1:B:210:ILE:HA	1.99	0.40
1:B:327:VAL:H	1:B:531:THR:CG2	2.35	0.40
1:B:362:VAL:HG13	1:B:525:CYS:O	2.21	0.40
1:C:405:ASN:N	1:C:504:GLY:O	2.54	0.40
1:C:645:THR:HG23	1:C:647:ALA:H	1.87	0.40
4:E:423:LEU:HD23	4:E:423:LEU:HA	1.88	0.40
4:F:48:TRP:HZ3	4:F:359:LEU:HB2	1.82	0.40
3:K:29:ILE:HB	3:K:71:THR:HA	2.03	0.40
1:A:1077:THR:OG1	1:A:1078:ALA:N	2.55	0.40
1:C:490:PHE:CE2	1:C:492:LEU:HB2	2.56	0.40
4:E:455:MET:HB2	4:E:455:MET:HE2	2.01	0.40
4:F:134:ASN:HB2	4:F:140:GLU:HB3	2.04	0.40
4:F:237:TYR:CZ	4:F:451:PRO:HG3	2.56	0.40
4:D:100:LEU:HD13	4:D:391:LEU:HD11	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1039/1247 (83%)	966 (93%)	72 (7%)	1 (0%)	48	80
1	B	1039/1247 (83%)	985 (95%)	54 (5%)	0	100	100
1	C	1039/1247 (83%)	974 (94%)	65 (6%)	0	100	100
2	G	222/228 (97%)	211 (95%)	11 (5%)	0	100	100
2	H	222/228 (97%)	213 (96%)	9 (4%)	0	100	100
2	I	222/228 (97%)	209 (94%)	13 (6%)	0	100	100
3	J	211/217 (97%)	202 (96%)	9 (4%)	0	100	100
3	K	211/217 (97%)	200 (95%)	11 (5%)	0	100	100
3	L	211/217 (97%)	202 (96%)	9 (4%)	0	100	100
4	D	594/602 (99%)	578 (97%)	16 (3%)	0	100	100
4	E	594/602 (99%)	579 (98%)	15 (2%)	0	100	100
4	F	594/602 (99%)	572 (96%)	22 (4%)	0	100	100
All	All	6198/6882 (90%)	5891 (95%)	306 (5%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	527	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	917/1083 (85%)	915 (100%)	2 (0%)	87	87
1	B	918/1083 (85%)	915 (100%)	3 (0%)	86	85
1	C	918/1083 (85%)	918 (100%)	0	100	100
2	G	186/191 (97%)	186 (100%)	0	100	100
2	H	186/191 (97%)	186 (100%)	0	100	100
2	I	186/191 (97%)	186 (100%)	0	100	100
3	J	177/183 (97%)	177 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	K	177/183 (97%)	177 (100%)	0	100	100
3	L	177/183 (97%)	177 (100%)	0	100	100
4	D	526/532 (99%)	526 (100%)	0	100	100
4	E	526/532 (99%)	526 (100%)	0	100	100
4	F	526/532 (99%)	525 (100%)	1 (0%)	87	87
All	All	5420/5967 (91%)	5414 (100%)	6 (0%)	87	89

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	525	CYS
1	A	534	VAL
1	B	331	ASN
1	B	525	CYS
1	B	529	LYS
4	F	360	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (57) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	ASN
1	A	414	GLN
1	A	474	GLN
1	A	762	GLN
1	A	804	GLN
1	A	895	GLN
1	A	935	GLN
1	A	1108	ASN
1	B	61	ASN
1	B	165	ASN
1	B	188	ASN
1	B	317	ASN
1	B	394	ASN
1	B	493	GLN
1	B	505	HIS
1	B	675	GLN
1	B	955	ASN
1	B	1011	GLN
1	C	188	ASN

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Mol	Chain	Res	Type
1	C	218	GLN
1	C	317	ASN
1	C	450	ASN
1	C	856	ASN
2	H	39	GLN
3	L	40	GLN
3	L	55	ASN
3	L	91	GLN
4	E	58	ASN
4	E	96	GLN
4	E	300	GLN
4	E	330	ASN
4	E	508	ASN
4	E	552	GLN
4	E	586	ASN
2	G	31	ASN
2	G	32	HIS
2	G	62	GLN
3	J	81	GLN
3	J	172	GLN
4	F	60	GLN
4	F	117	ASN
4	F	121	ASN
4	F	194	ASN
4	F	300	GLN
4	F	330	ASN
4	F	586	ASN
2	I	3	GLN
2	I	39	GLN
3	K	40	GLN
3	K	199	GLN
4	D	96	GLN
4	D	221	GLN
4	D	373	HIS
4	D	388	GLN
4	D	508	ASN
4	D	524	GLN
4	D	586	ASN

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 ⓘ

12 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	M	1	5,1	14,14,15	0.28	0	17,19,21	0.42	0
5	NAG	M	2	5	14,14,15	0.21	0	17,19,21	0.47	0
5	NAG	N	1	5,1	14,14,15	0.24	0	17,19,21	0.43	0
5	NAG	N	2	5	14,14,15	0.24	0	17,19,21	0.43	0
5	NAG	O	1	5,1	14,14,15	0.20	0	17,19,21	0.42	0
5	NAG	O	2	5	14,14,15	0.24	0	17,19,21	0.50	0
5	NAG	P	1	5,1	14,14,15	0.27	0	17,19,21	0.43	0
5	NAG	P	2	5	14,14,15	0.23	0	17,19,21	0.45	0
5	NAG	Q	1	5,1	14,14,15	0.22	0	17,19,21	0.48	0
5	NAG	Q	2	5	14,14,15	0.22	0	17,19,21	0.46	0
5	NAG	R	1	5,1	14,14,15	0.31	0	17,19,21	0.41	0
5	NAG	R	2	5	14,14,15	0.23	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	M	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	M	2	5	-	2/6/23/26	0/1/1/1
5	NAG	N	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	N	2	5	-	2/6/23/26	0/1/1/1
5	NAG	O	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	O	2	5	-	2/6/23/26	0/1/1/1
5	NAG	P	1	5,1	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	P	2	5	-	2/6/23/26	0/1/1/1
5	NAG	Q	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	Q	2	5	-	2/6/23/26	0/1/1/1
5	NAG	R	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	R	2	5	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (19) torsion outliers are listed below:

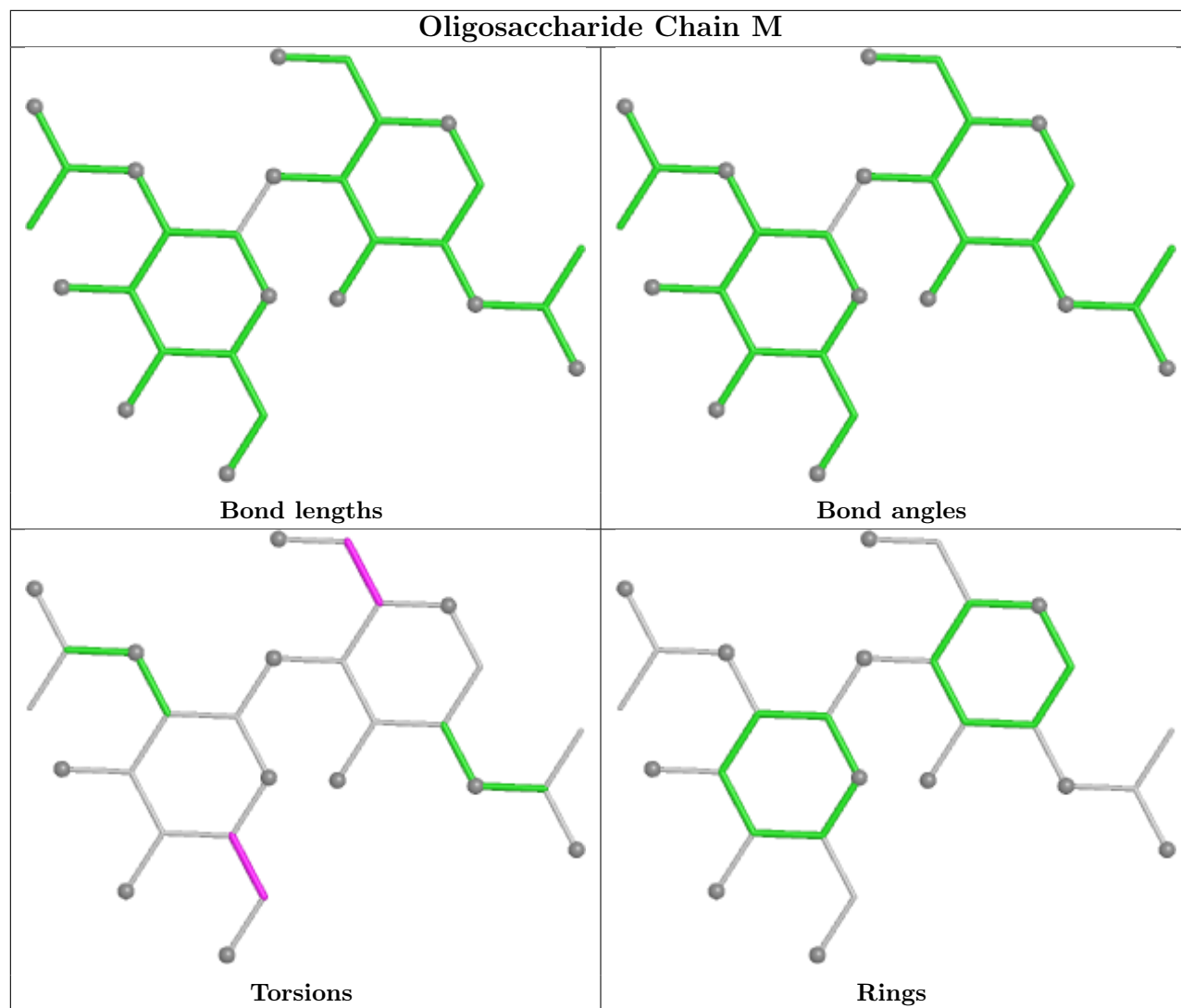
Mol	Chain	Res	Type	Atoms
5	R	2	NAG	O5-C5-C6-O6
5	N	2	NAG	O5-C5-C6-O6
5	O	2	NAG	O5-C5-C6-O6
5	M	2	NAG	O5-C5-C6-O6
5	R	2	NAG	C4-C5-C6-O6
5	N	1	NAG	O5-C5-C6-O6
5	N	2	NAG	C4-C5-C6-O6
5	M	2	NAG	C4-C5-C6-O6
5	N	1	NAG	C4-C5-C6-O6
5	O	2	NAG	C4-C5-C6-O6
5	P	2	NAG	O5-C5-C6-O6
5	P	2	NAG	C4-C5-C6-O6
5	R	1	NAG	O5-C5-C6-O6
5	P	1	NAG	O5-C5-C6-O6
5	Q	2	NAG	C4-C5-C6-O6
5	Q	2	NAG	O5-C5-C6-O6
5	M	1	NAG	C4-C5-C6-O6
5	O	1	NAG	C4-C5-C6-O6
5	M	1	NAG	O5-C5-C6-O6

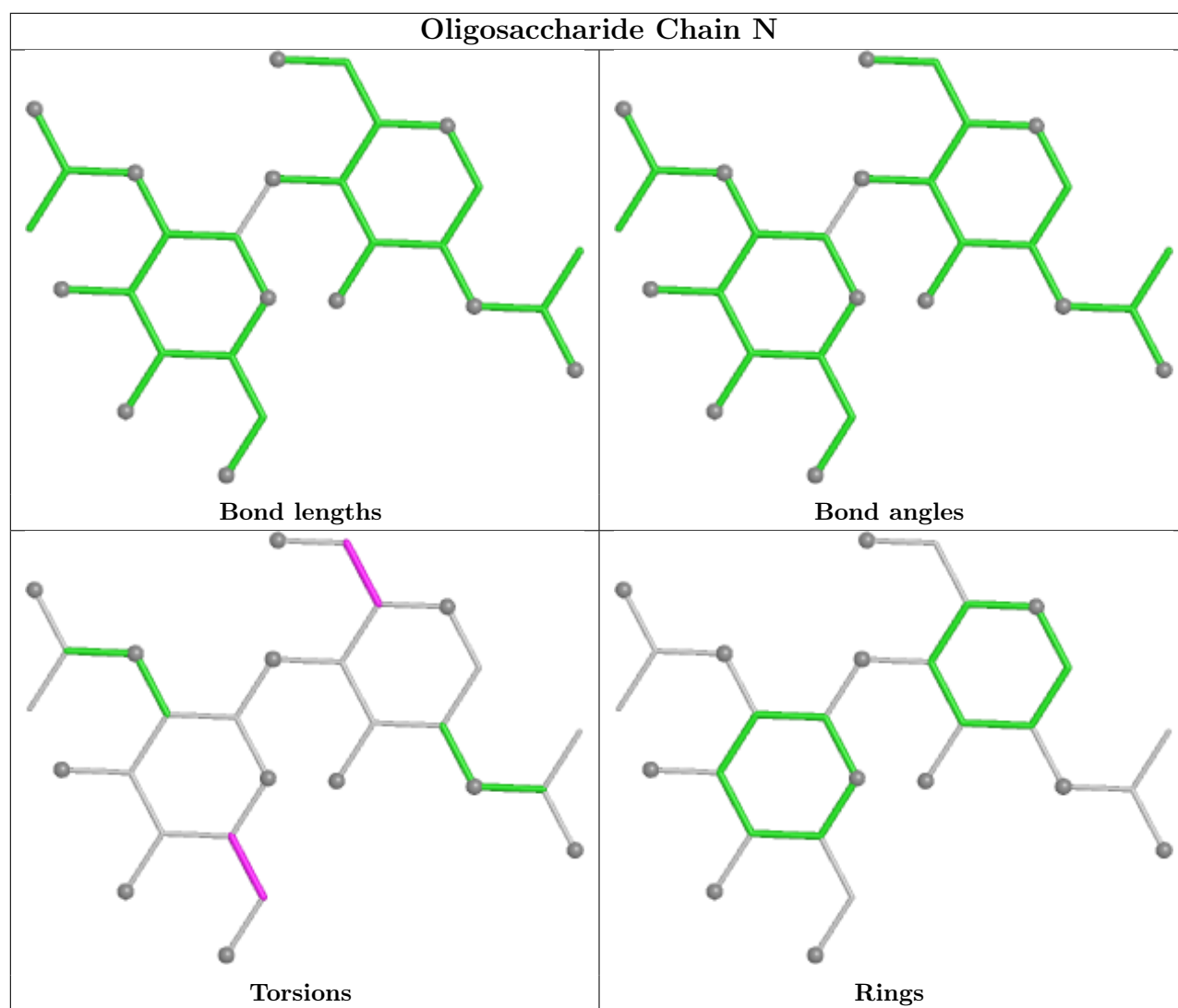
There are no ring outliers.

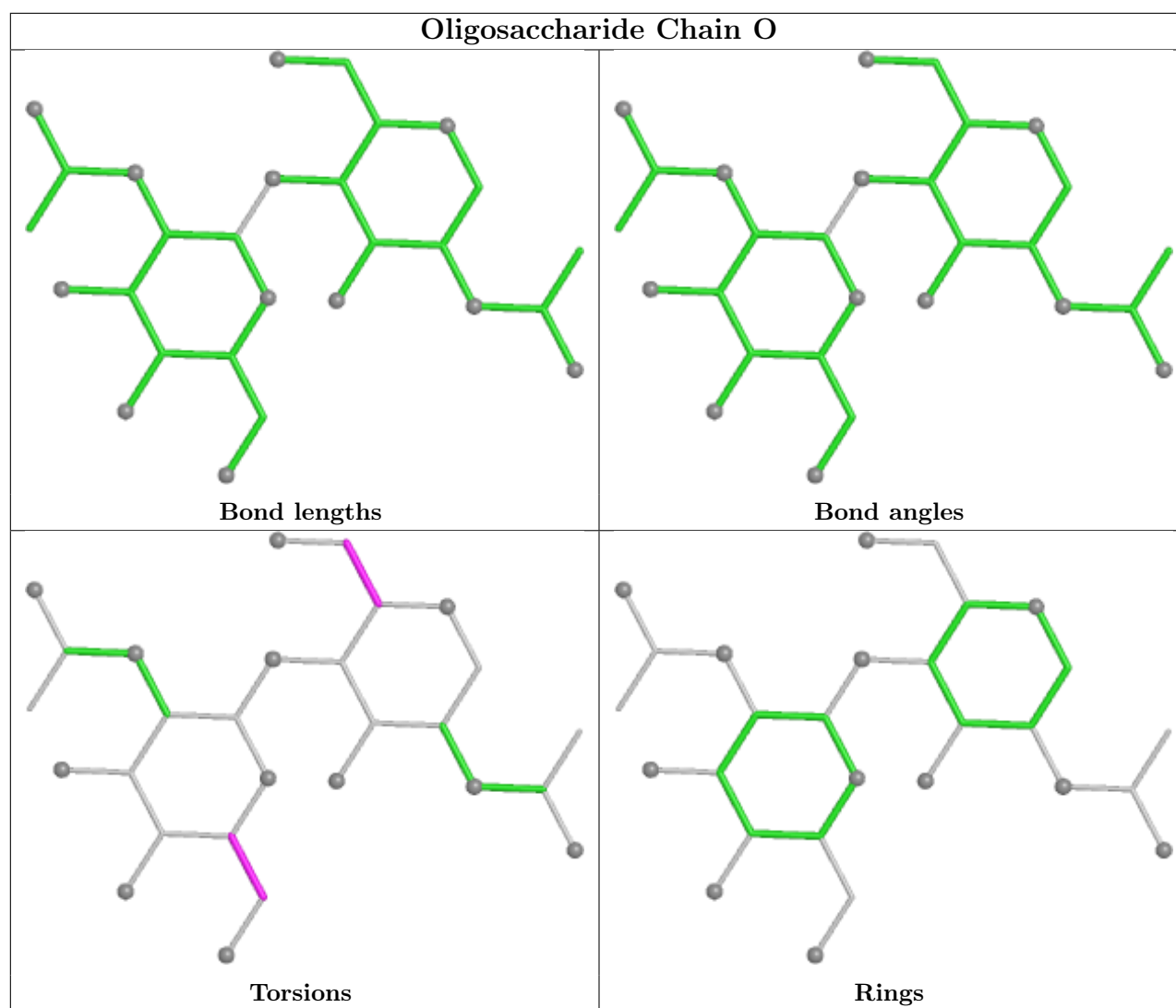
1 monomer is involved in 1 short contact:

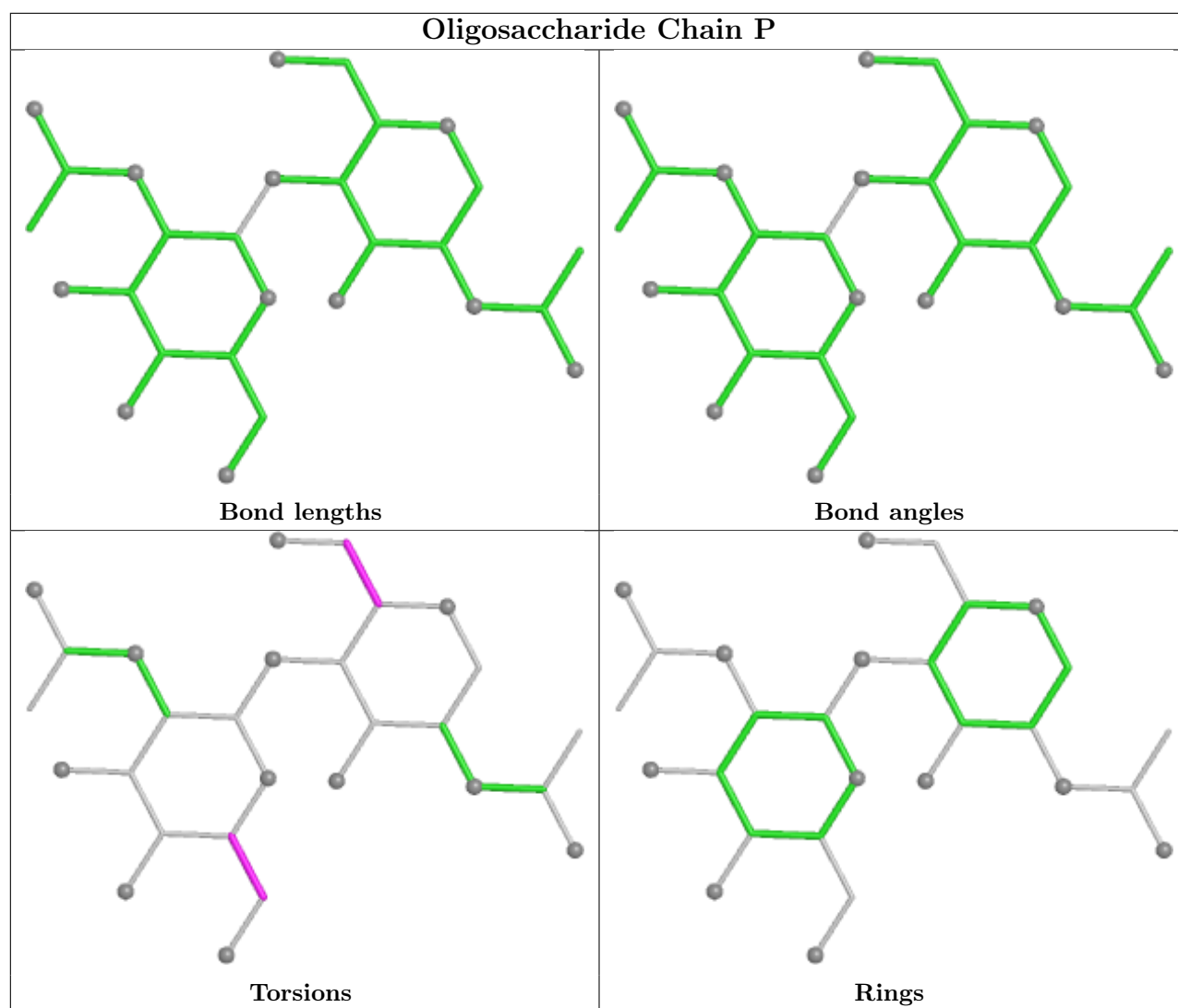
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	M	1	NAG	1	0

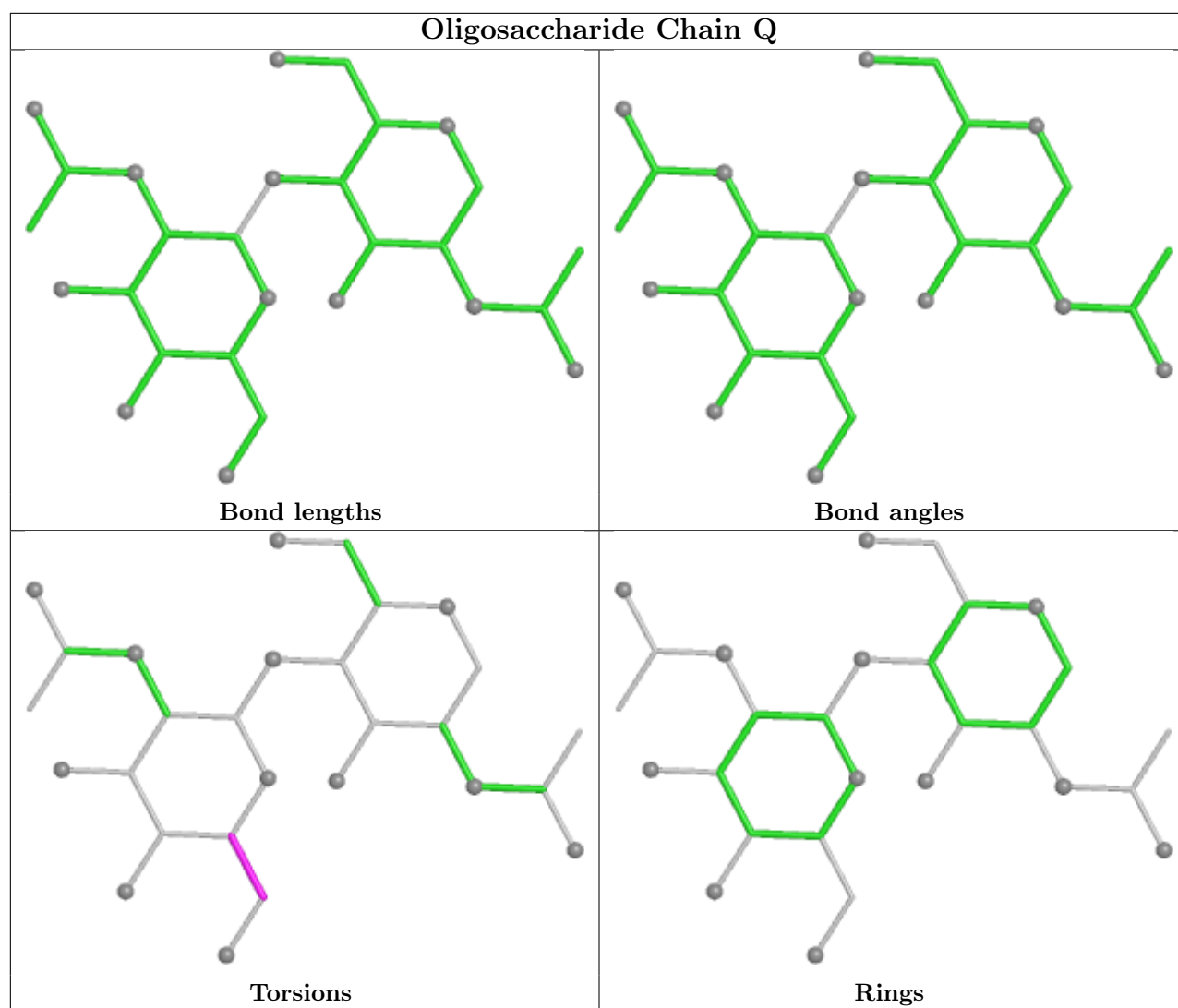
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

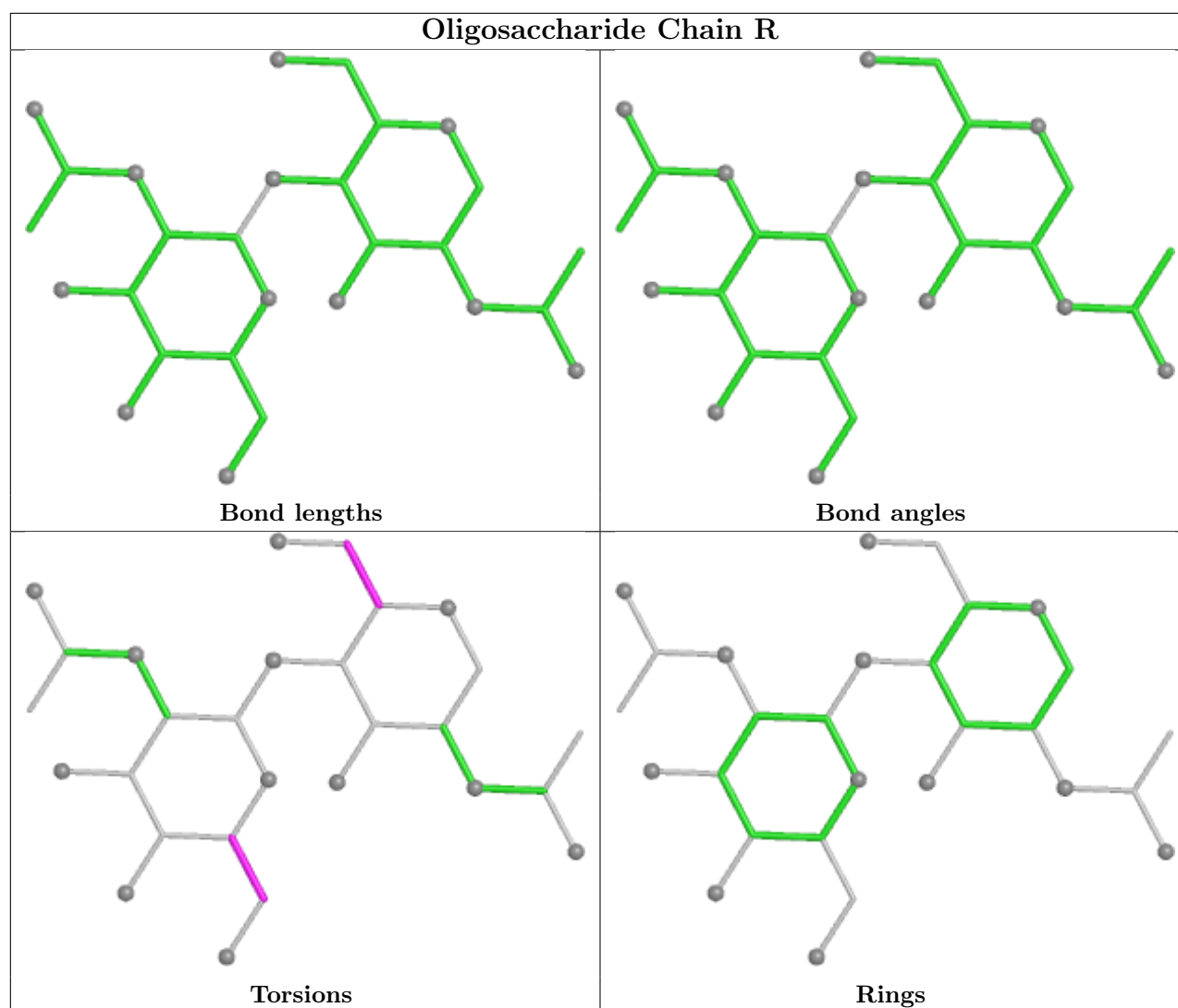












5.6 Ligand geometry [i](#)

Of 39 ligands modelled in this entry, 3 are monoatomic - leaving 36 for Mogul analysis.

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	$\# Z > 2$
6	NAG	C	1306	1	14,14,15	0.21	0	17,19,21	0.49	0
6	NAG	E	903	4	14,14,15	0.34	0	17,19,21	0.64	0
6	NAG	C	1304	1	14,14,15	0.27	0	17,19,21	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	E	904	4	14,14,15	0.30	0	17,19,21	0.40	0
6	NAG	F	902	4	14,14,15	0.30	0	17,19,21	0.43	0
6	NAG	A	1302	1	14,14,15	0.23	0	17,19,21	0.41	0
6	NAG	B	1302	1	14,14,15	0.29	0	17,19,21	0.88	0
6	NAG	A	1303	1	14,14,15	0.50	0	17,19,21	0.48	0
6	NAG	D	902	4	14,14,15	0.24	0	17,19,21	0.43	0
6	NAG	A	1307	1	14,14,15	0.34	0	17,19,21	0.36	0
6	NAG	C	1302	1	14,14,15	0.26	0	17,19,21	0.38	0
6	NAG	B	1308	1	14,14,15	0.25	0	17,19,21	0.40	0
6	NAG	E	905	4	14,14,15	0.28	0	17,19,21	0.46	0
6	NAG	D	903	4	14,14,15	0.34	0	17,19,21	0.72	0
6	NAG	D	904	4	14,14,15	0.27	0	17,19,21	0.42	0
6	NAG	A	1305	1	14,14,15	0.21	0	17,19,21	0.39	0
6	NAG	A	1308	1	14,14,15	0.31	0	17,19,21	0.42	0
6	NAG	B	1305	1	14,14,15	0.24	0	17,19,21	0.45	0
6	NAG	B	1306	1	14,14,15	0.21	0	17,19,21	0.54	0
6	NAG	C	1303	1	14,14,15	0.38	0	17,19,21	0.43	0
6	NAG	C	1307	1	14,14,15	0.37	0	17,19,21	0.37	0
6	NAG	F	904	4	14,14,15	0.24	0	17,19,21	0.44	0
6	NAG	F	903	4	14,14,15	0.32	0	17,19,21	0.54	0
6	NAG	C	1301	1	14,14,15	0.27	0	17,19,21	0.72	0
6	NAG	B	1303	1	14,14,15	0.65	1 (7%)	17,19,21	0.62	1 (5%)
6	NAG	A	1304	1	14,14,15	0.22	0	17,19,21	0.51	0
6	NAG	B	1304	1	14,14,15	0.30	0	17,19,21	0.48	0
6	NAG	F	905	4	14,14,15	0.25	0	17,19,21	0.44	0
6	NAG	A	1306	1	14,14,15	0.23	0	17,19,21	0.50	0
6	NAG	E	902	4	14,14,15	0.26	0	17,19,21	0.46	0
6	NAG	B	1307	1	14,14,15	0.41	0	17,19,21	0.41	0
6	NAG	B	1301	1	14,14,15	0.34	0	17,19,21	0.52	0
6	NAG	A	1301	1	14,14,15	0.23	0	17,19,21	0.45	0
6	NAG	C	1305	1	14,14,15	0.25	0	17,19,21	0.45	0
6	NAG	C	1308	1	14,14,15	0.24	0	17,19,21	0.40	0
6	NAG	D	905	4	14,14,15	0.23	0	17,19,21	0.44	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
6	NAG	C	1306	1	-	1/6/23/26	0/1/1/1
6	NAG	E	903	4	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	C	1304	1	-	2/6/23/26	0/1/1/1
6	NAG	E	904	4	-	2/6/23/26	0/1/1/1
6	NAG	F	902	4	-	2/6/23/26	0/1/1/1
6	NAG	A	1302	1	-	2/6/23/26	0/1/1/1
6	NAG	B	1302	1	-	4/6/23/26	0/1/1/1
6	NAG	A	1303	1	-	2/6/23/26	0/1/1/1
6	NAG	D	902	4	-	2/6/23/26	0/1/1/1
6	NAG	A	1307	1	-	2/6/23/26	0/1/1/1
6	NAG	C	1302	1	-	4/6/23/26	0/1/1/1
6	NAG	B	1308	1	-	2/6/23/26	0/1/1/1
6	NAG	E	905	4	-	0/6/23/26	0/1/1/1
6	NAG	D	903	4	-	2/6/23/26	0/1/1/1
6	NAG	D	904	4	-	0/6/23/26	0/1/1/1
6	NAG	A	1305	1	-	2/6/23/26	0/1/1/1
6	NAG	A	1308	1	-	1/6/23/26	0/1/1/1
6	NAG	B	1305	1	-	0/6/23/26	0/1/1/1
6	NAG	B	1306	1	-	2/6/23/26	0/1/1/1
6	NAG	C	1303	1	-	2/6/23/26	0/1/1/1
6	NAG	C	1307	1	-	2/6/23/26	0/1/1/1
6	NAG	F	904	4	-	0/6/23/26	0/1/1/1
6	NAG	F	903	4	-	3/6/23/26	0/1/1/1
6	NAG	C	1301	1	-	3/6/23/26	0/1/1/1
6	NAG	B	1303	1	-	2/6/23/26	0/1/1/1
6	NAG	A	1304	1	-	2/6/23/26	0/1/1/1
6	NAG	B	1304	1	-	2/6/23/26	0/1/1/1
6	NAG	F	905	4	-	2/6/23/26	0/1/1/1
6	NAG	A	1306	1	-	0/6/23/26	0/1/1/1
6	NAG	E	902	4	-	2/6/23/26	0/1/1/1
6	NAG	B	1307	1	-	3/6/23/26	0/1/1/1
6	NAG	B	1301	1	-	0/6/23/26	0/1/1/1
6	NAG	A	1301	1	-	0/6/23/26	0/1/1/1
6	NAG	C	1305	1	-	2/6/23/26	0/1/1/1
6	NAG	C	1308	1	-	0/6/23/26	0/1/1/1
6	NAG	D	905	4	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	1303	NAG	O5-C1	2.01	1.46	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	1303	NAG	C1-O5-C5	2.22	115.20	112.19

There are no chirality outliers.

All (61) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	1302	NAG	C8-C7-N2-C2
6	B	1302	NAG	O7-C7-N2-C2
6	C	1301	NAG	C3-C2-N2-C7
6	C	1301	NAG	C8-C7-N2-C2
6	C	1301	NAG	O7-C7-N2-C2
6	F	903	NAG	O7-C7-N2-C2
6	A	1307	NAG	C4-C5-C6-O6
6	C	1302	NAG	O5-C5-C6-O6
6	F	903	NAG	C8-C7-N2-C2
6	F	905	NAG	O5-C5-C6-O6
6	A	1307	NAG	O5-C5-C6-O6
6	B	1307	NAG	C4-C5-C6-O6
6	B	1306	NAG	C4-C5-C6-O6
6	C	1305	NAG	C4-C5-C6-O6
6	E	902	NAG	C4-C5-C6-O6
6	A	1304	NAG	O5-C5-C6-O6
6	B	1306	NAG	O5-C5-C6-O6
6	B	1304	NAG	O5-C5-C6-O6
6	D	902	NAG	O5-C5-C6-O6
6	D	905	NAG	O5-C5-C6-O6
6	C	1302	NAG	C4-C5-C6-O6
6	A	1303	NAG	O5-C5-C6-O6
6	B	1307	NAG	O5-C5-C6-O6
6	E	902	NAG	O5-C5-C6-O6
6	E	904	NAG	O5-C5-C6-O6
6	C	1307	NAG	C4-C5-C6-O6
6	F	905	NAG	C4-C5-C6-O6
6	D	903	NAG	C8-C7-N2-C2
6	B	1303	NAG	O5-C5-C6-O6
6	B	1302	NAG	C1-C2-N2-C7
6	A	1305	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
6	B	1304	NAG	C4-C5-C6-O6
6	D	905	NAG	C4-C5-C6-O6
6	E	904	NAG	C4-C5-C6-O6
6	B	1303	NAG	C4-C5-C6-O6
6	A	1302	NAG	C8-C7-N2-C2
6	A	1302	NAG	O7-C7-N2-C2
6	C	1302	NAG	C8-C7-N2-C2
6	C	1302	NAG	O7-C7-N2-C2
6	E	903	NAG	C8-C7-N2-C2
6	D	903	NAG	O7-C7-N2-C2
6	B	1308	NAG	O5-C5-C6-O6
6	A	1303	NAG	C4-C5-C6-O6
6	A	1304	NAG	C4-C5-C6-O6
6	C	1307	NAG	O5-C5-C6-O6
6	C	1303	NAG	C4-C5-C6-O6
6	F	902	NAG	O5-C5-C6-O6
6	C	1304	NAG	O5-C5-C6-O6
6	C	1305	NAG	O5-C5-C6-O6
6	F	903	NAG	C1-C2-N2-C7
6	B	1308	NAG	C4-C5-C6-O6
6	E	903	NAG	O7-C7-N2-C2
6	D	902	NAG	C4-C5-C6-O6
6	C	1303	NAG	O5-C5-C6-O6
6	C	1304	NAG	C4-C5-C6-O6
6	A	1305	NAG	C4-C5-C6-O6
6	B	1302	NAG	C3-C2-N2-C7
6	C	1306	NAG	C4-C5-C6-O6
6	B	1307	NAG	C1-C2-N2-C7
6	A	1308	NAG	O5-C5-C6-O6
6	F	902	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1302	NAG	2	0
6	C	1301	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

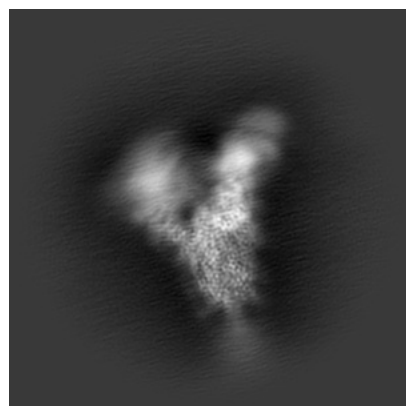
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-66892. These allow visual inspection of the internal detail of the map and identification of artifacts.

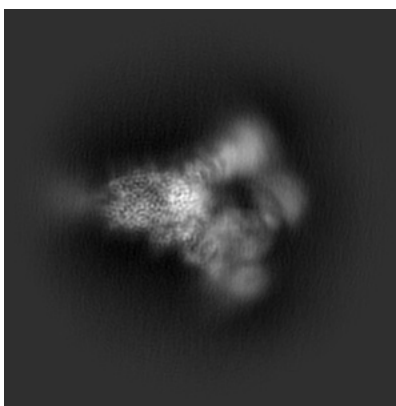
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

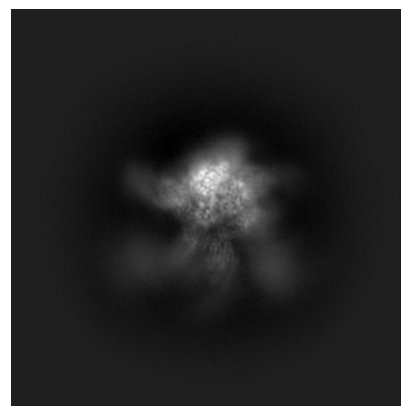
6.1.1 Primary map



X

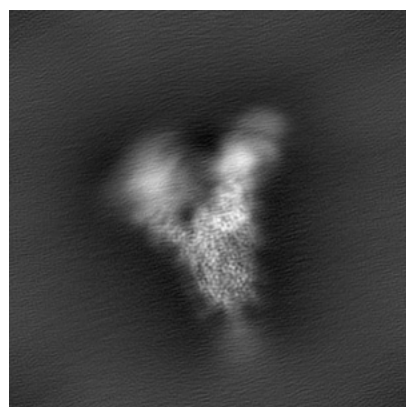


Y

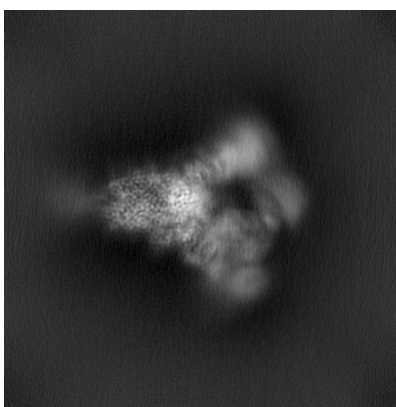


Z

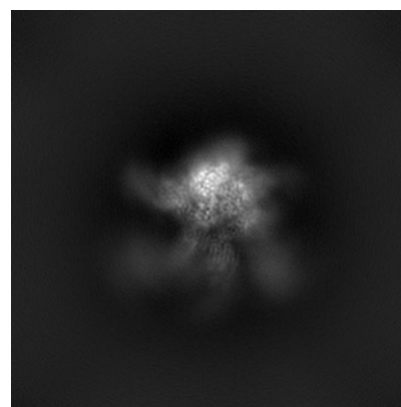
6.1.2 Raw map



X



Y

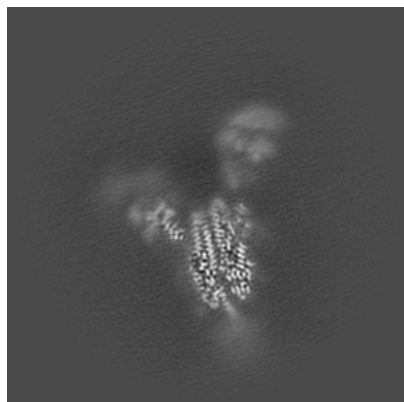


Z

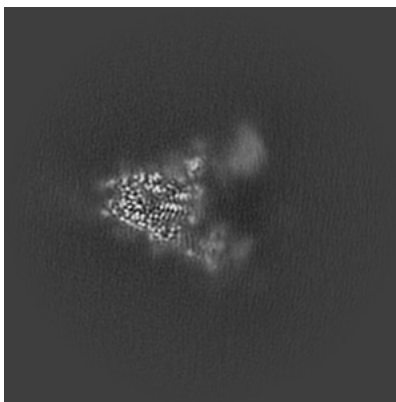
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

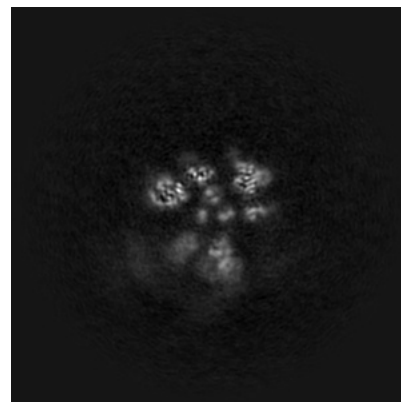
6.2.1 Primary map



X Index: 128

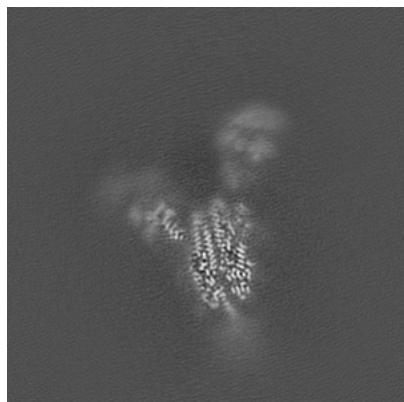


Y Index: 128

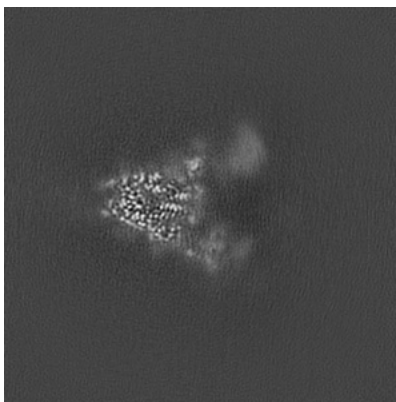


Z Index: 128

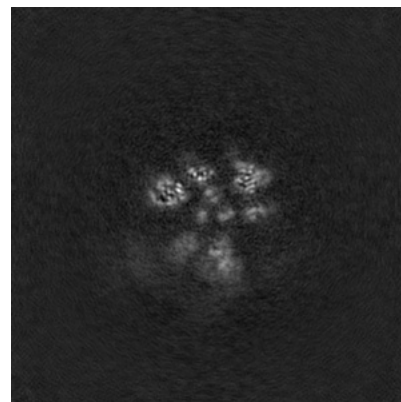
6.2.2 Raw map



X Index: 128



Y Index: 128

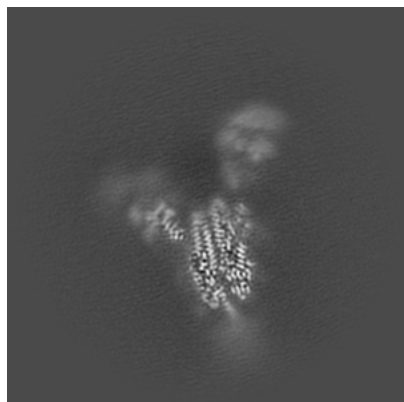


Z Index: 128

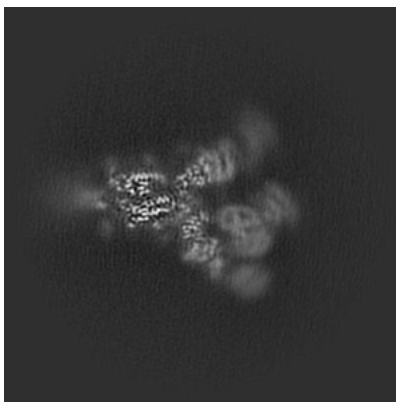
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

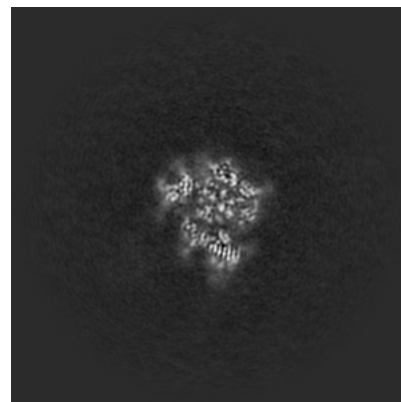
6.3.1 Primary map



X Index: 128

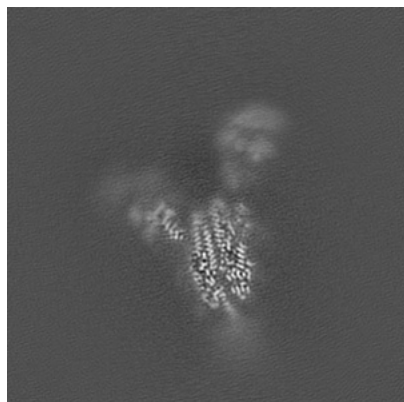


Y Index: 143

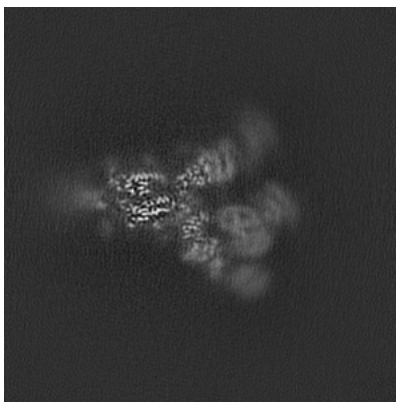


Z Index: 116

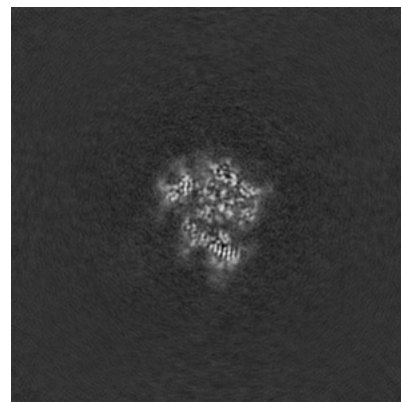
6.3.2 Raw map



X Index: 128



Y Index: 143

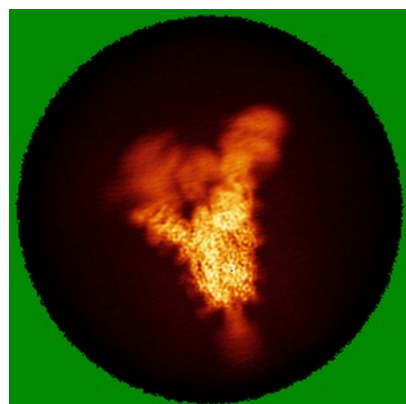


Z Index: 116

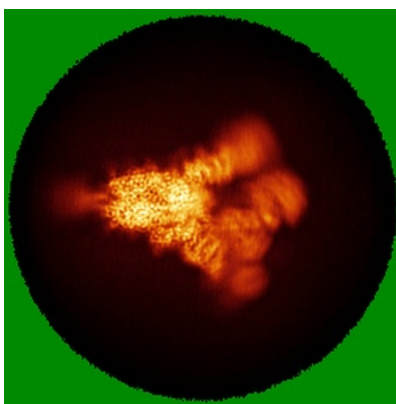
The images above show the largest variance slices of the map in three orthogonal directions.

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

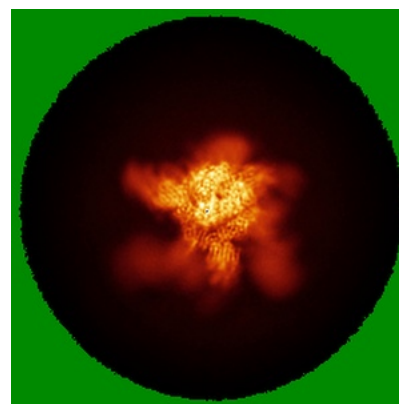
6.4.1 Primary map



X



Y

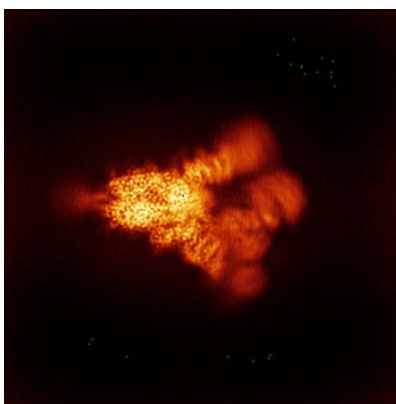


Z

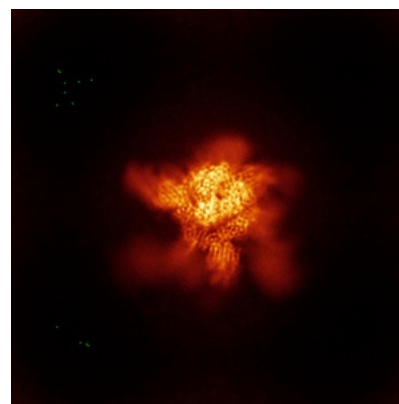
6.4.2 Raw map



X



Y

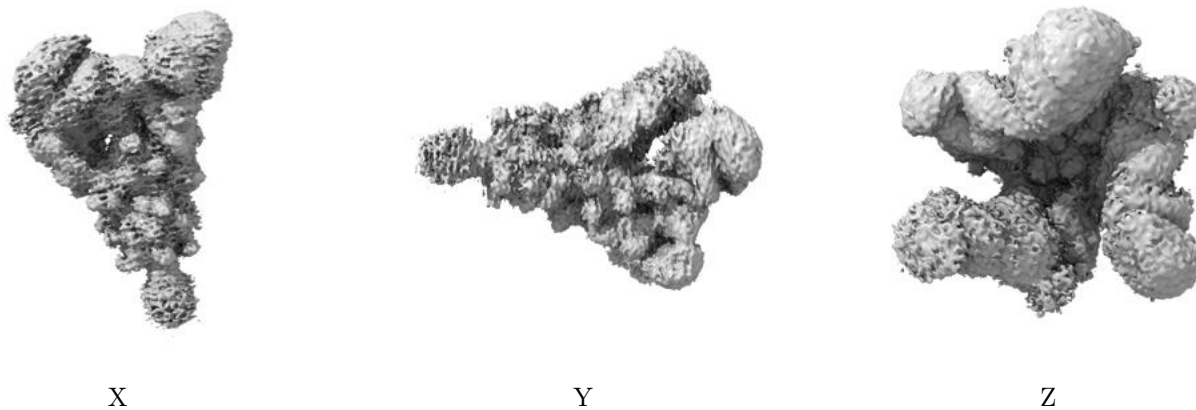


Z

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

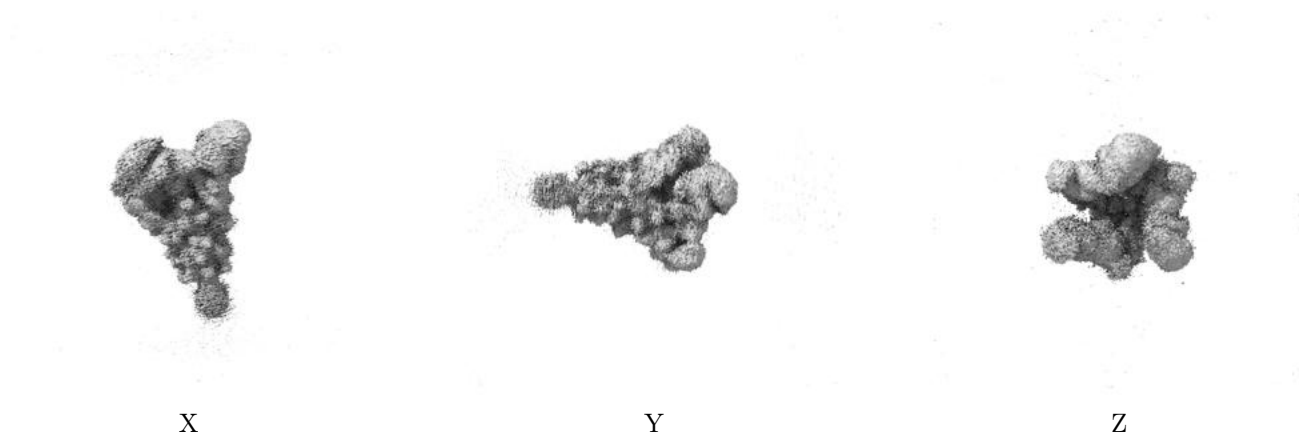
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.07. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

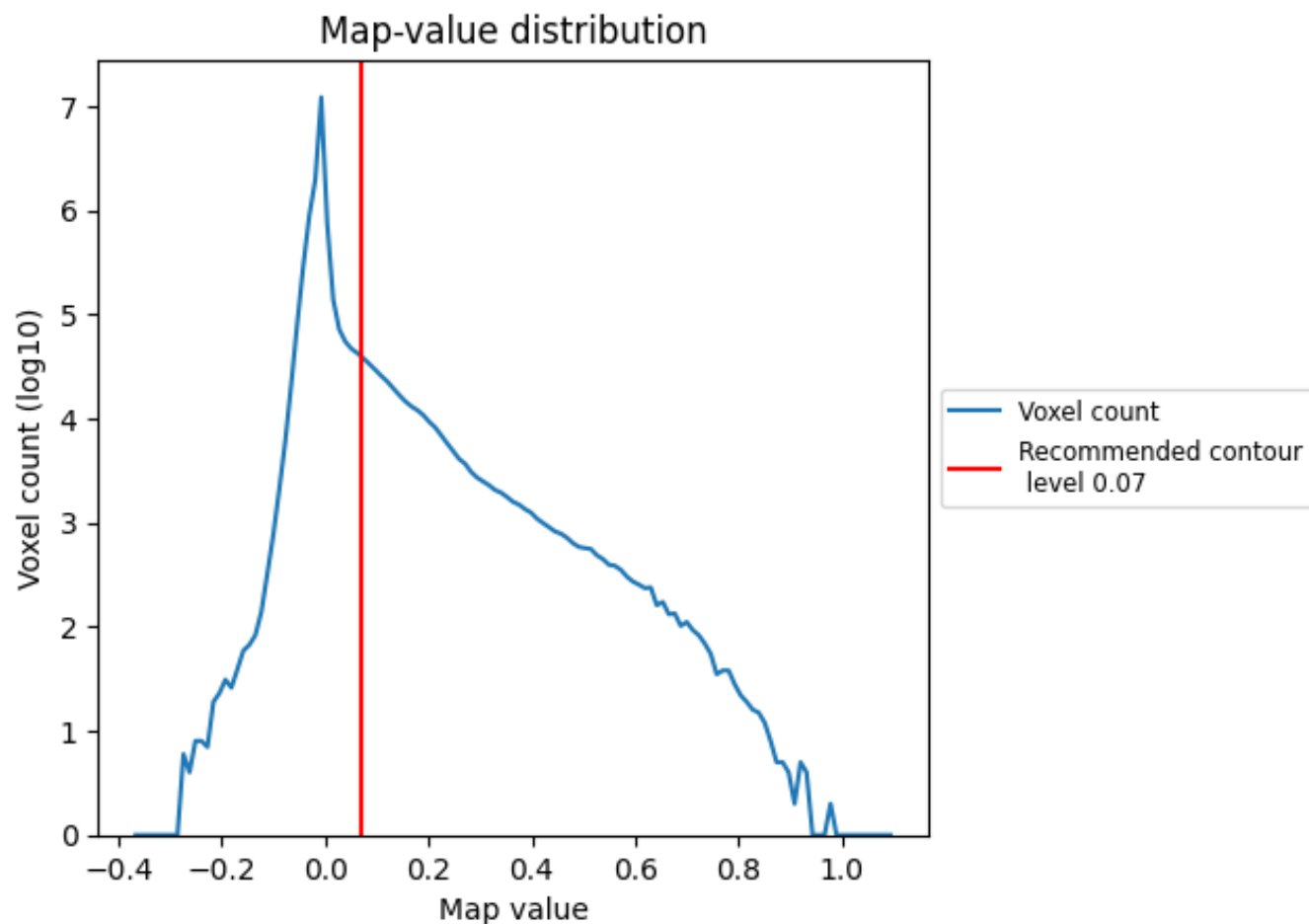
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

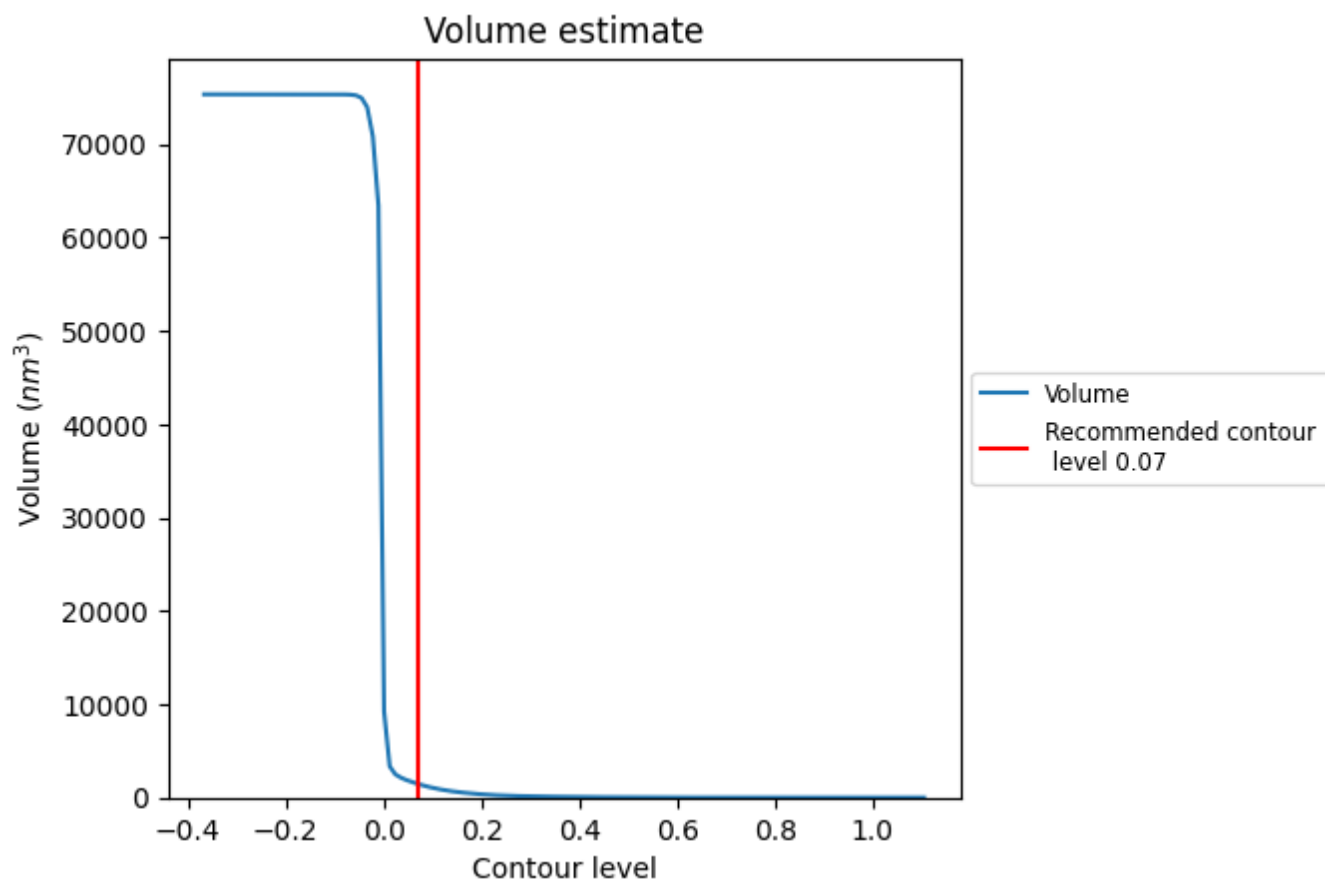
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

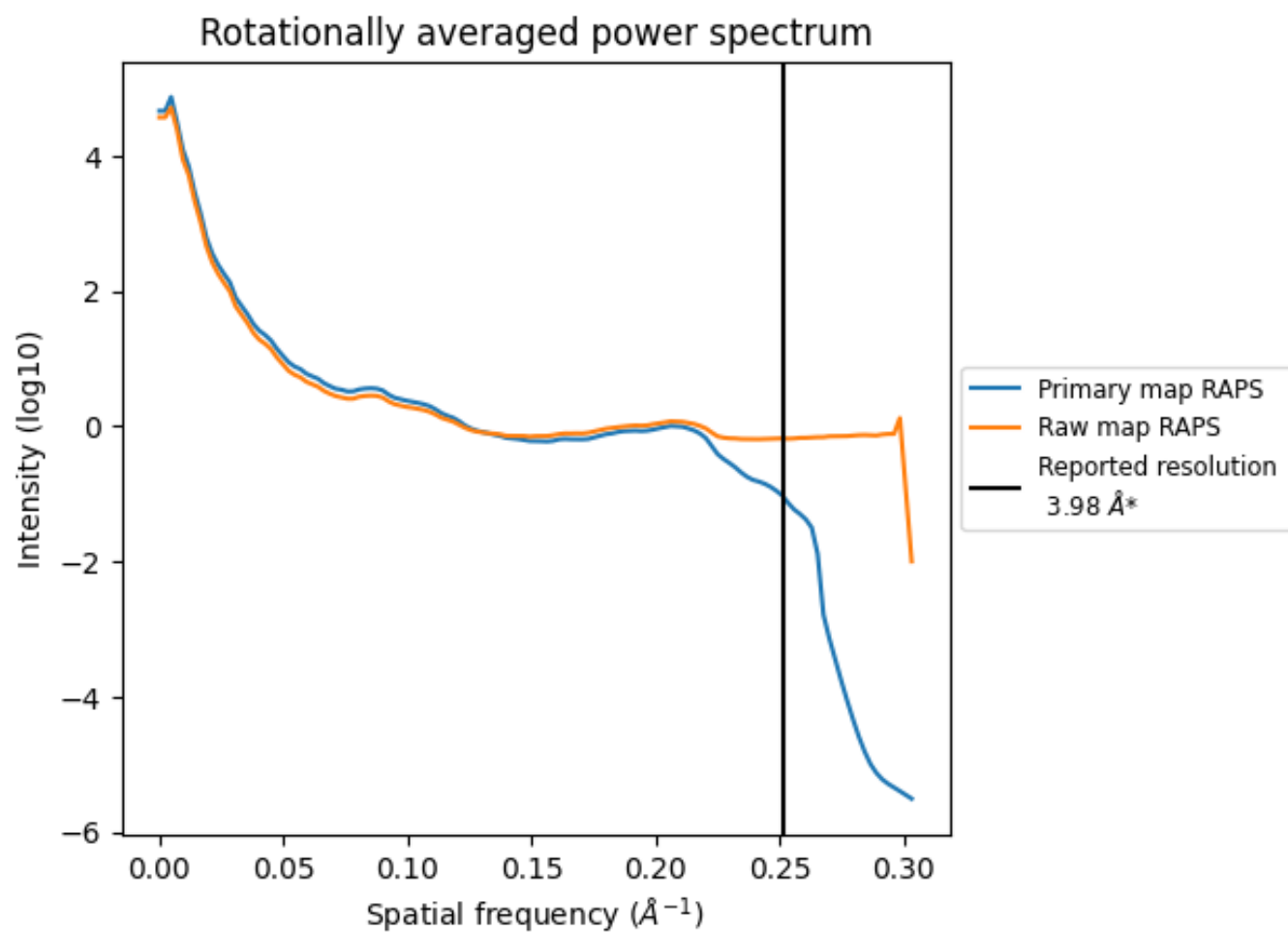
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1471 nm^3 ; this corresponds to an approximate mass of 1329 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

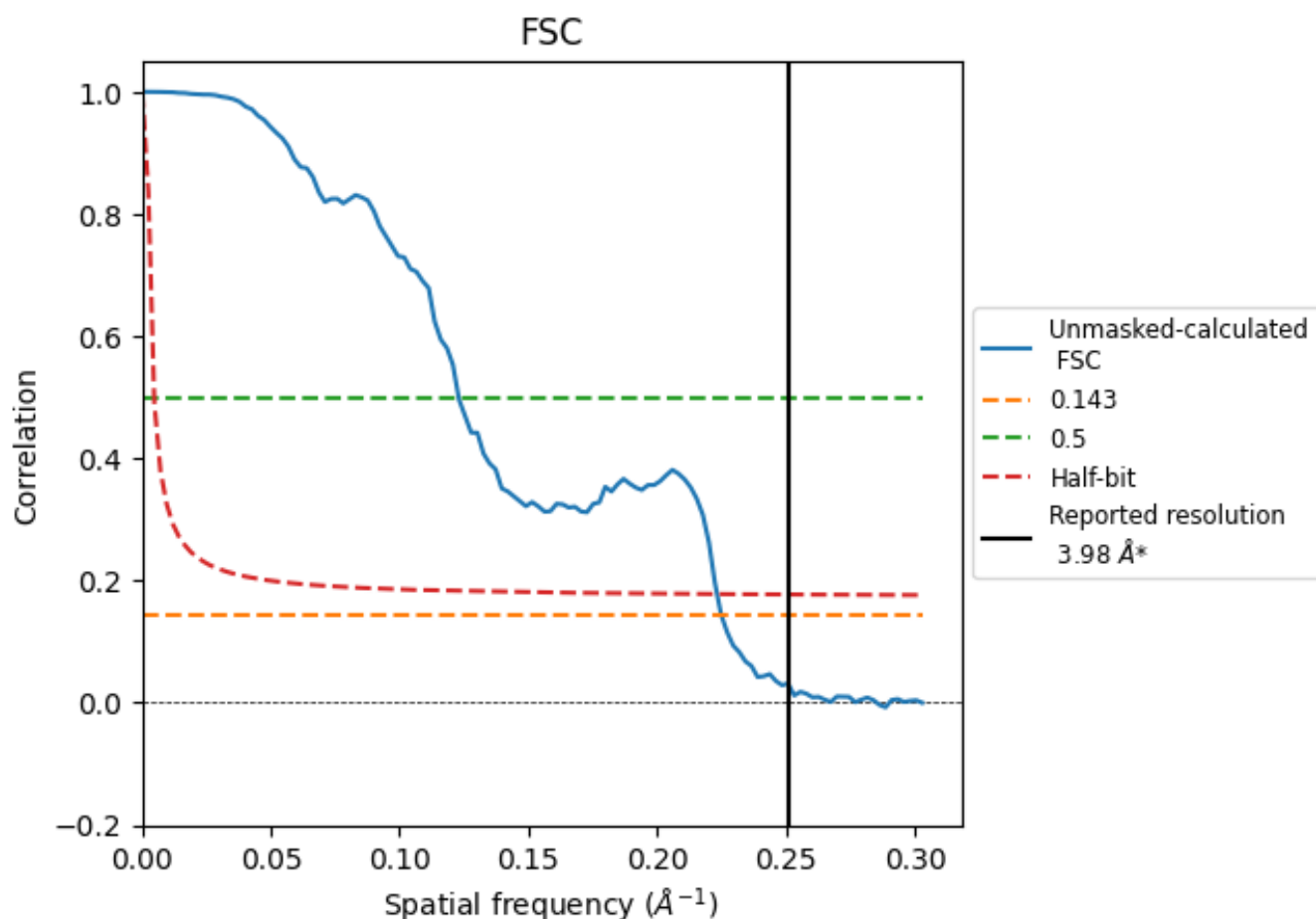


*Reported resolution corresponds to spatial frequency of 0.251 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.251 Å⁻¹

8.2 Resolution estimates [i](#)

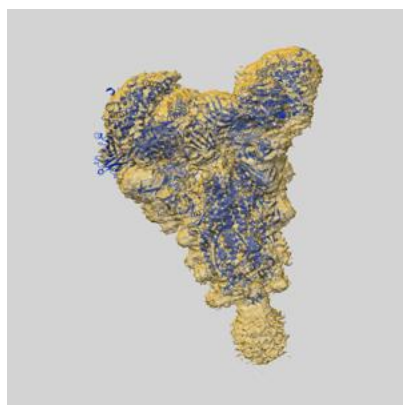
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.98	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.44	8.14	4.48

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.44 differs from the reported value 3.98 by more than 10 %

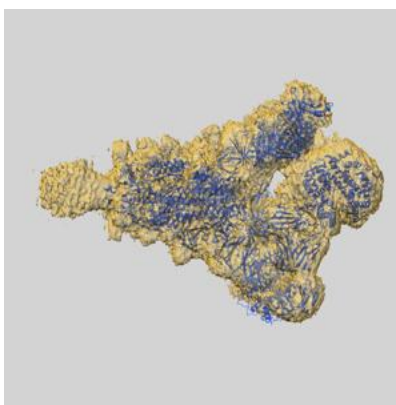
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-66892 and PDB model 9XI0. Per-residue inclusion information can be found in section 3 on page 15.

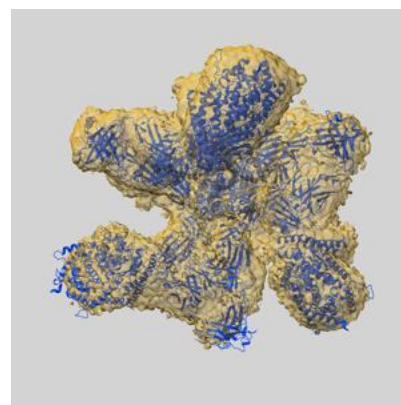
9.1 Map-model overlay [i](#)



X



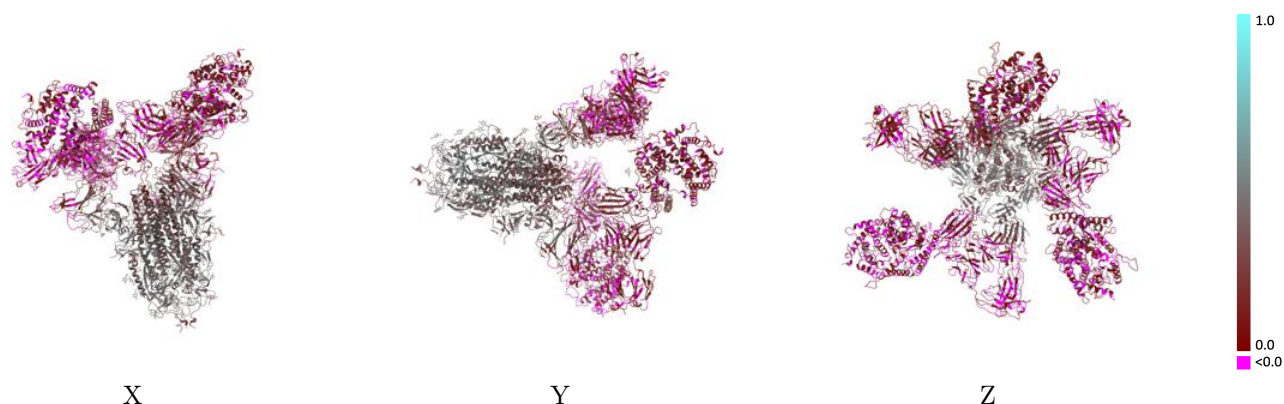
Y



Z

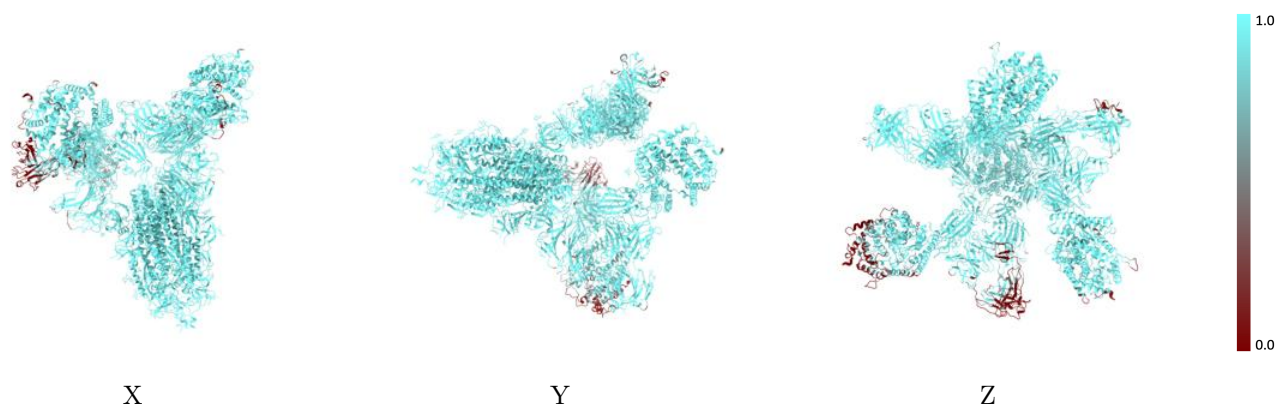
The images above show the 3D surface view of the map at the recommended contour level 0.07 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



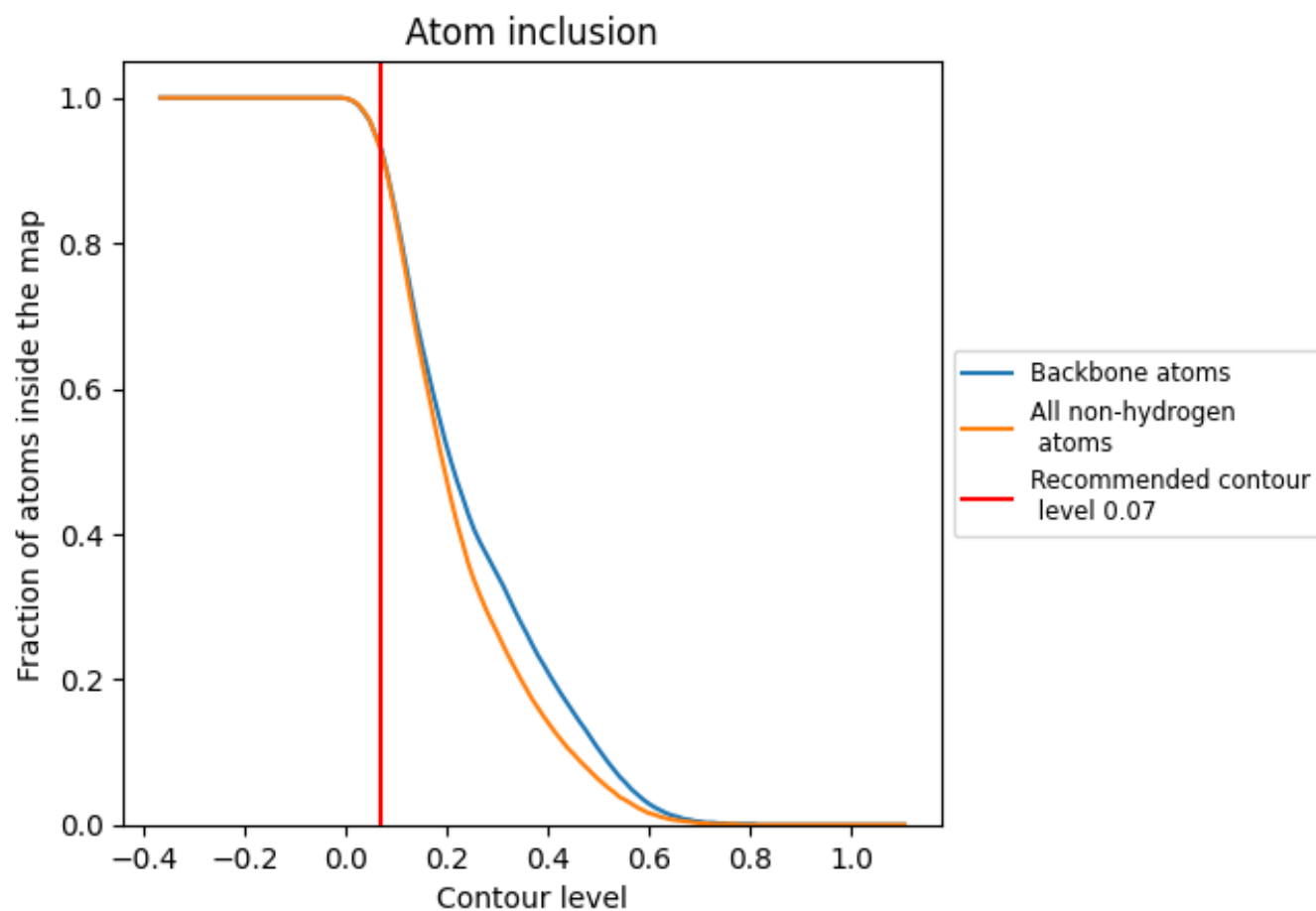
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.07).

9.4 Atom inclusion ⓘ



At the recommended contour level, 93% of all backbone atoms, 93% 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 (0.07) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9270	 0.2000
A	 0.9960	 0.3100
B	 0.9910	 0.3100
C	 0.9960	 0.3210
D	 0.7040	 0.0530
E	 0.9440	 0.0740
F	 0.9930	 0.1180
G	 0.9110	 0.0820
H	 0.6570	 0.0590
I	 0.9840	 0.1220
J	 0.8930	 0.0740
K	 0.9990	 0.1060
L	 0.4990	 0.0490
M	 1.0000	 0.4320
N	 1.0000	 0.3880
O	 1.0000	 0.4440
P	 1.0000	 0.4030
Q	 1.0000	 0.4370
R	 1.0000	 0.4080

