



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 04:08 AM UTC

PDB ID : 7WVQ / pdb_00007wvq
EMDB ID : EMD-32857
Title : Cryo-EM structure of SARS-CoV-2 Omicron Spike protein with human ACE2 receptor, C3 state
Authors : Han, W.Y.; Wang, Y.F.
Deposited on : 2022-02-10
Resolution : 4.04 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

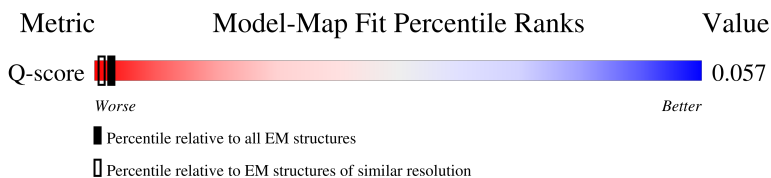
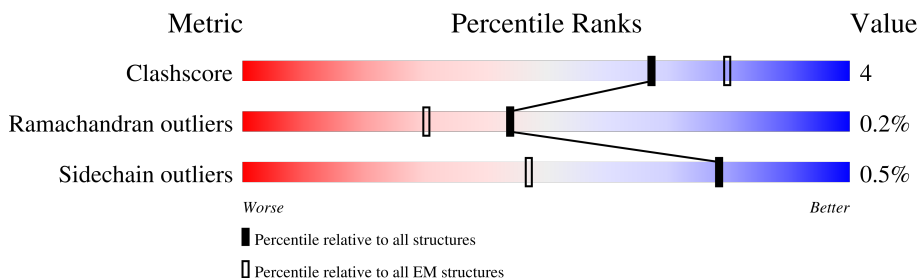
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 4.04 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	6625 (3.54 - 4.54)

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	B	1258	15% 73% 8% 17%
1	C	1258	19% 75% 7% 17%
1	D	1258	20% 77% 7% 17%
2	A	625	50% 87% 8% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 29527 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	D	1048	Total	C	N	O	S	0	0
			8219	5260	1369	1552	38		
1	C	1048	Total	C	N	O	S	0	0
			8219	5260	1369	1552	38		
1	B	1048	Total	C	N	O	S	0	0
			8219	5260	1369	1552	38		

There are 294 discrepancies between the modelled and reference sequences:

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

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Chain	Residue	Modelled	Actual	Comment	Reference
D	493	ARG	GLN	variant	UNP P0DTC2
D	496	SER	GLY	variant	UNP P0DTC2
D	498	ARG	GLN	variant	UNP P0DTC2
D	501	TYR	ASN	variant	UNP P0DTC2
D	505	HIS	TYR	variant	UNP P0DTC2
D	547	LYS	THR	variant	UNP P0DTC2
D	614	GLY	ASP	variant	UNP P0DTC2
D	655	TYR	HIS	variant	UNP P0DTC2
D	679	LYS	ASN	variant	UNP P0DTC2
D	681	HIS	PRO	variant	UNP P0DTC2
D	682	GLY	ARG	variant	UNP P0DTC2
D	683	SER	ARG	variant	UNP P0DTC2
D	685	SER	ARG	variant	UNP P0DTC2
D	764	LYS	ASN	variant	UNP P0DTC2
D	796	TYR	ASP	variant	UNP P0DTC2
D	856	LYS	ASN	variant	UNP P0DTC2
D	954	HIS	GLN	variant	UNP P0DTC2
D	969	LYS	ASN	variant	UNP P0DTC2
D	981	PHE	LEU	variant	UNP P0DTC2
D	986	PRO	LYS	variant	UNP P0DTC2
D	987	PRO	VAL	variant	UNP P0DTC2
D	1209	GLY	-	expression tag	UNP P0DTC2
D	1210	SER	-	expression tag	UNP P0DTC2
D	1211	GLY	-	expression tag	UNP P0DTC2
D	1212	TYR	-	expression tag	UNP P0DTC2
D	1213	ILE	-	expression tag	UNP P0DTC2
D	1214	PRO	-	expression tag	UNP P0DTC2
D	1215	GLU	-	expression tag	UNP P0DTC2
D	1216	ALA	-	expression tag	UNP P0DTC2
D	1217	PRO	-	expression tag	UNP P0DTC2
D	1218	ARG	-	expression tag	UNP P0DTC2
D	1219	ASP	-	expression tag	UNP P0DTC2
D	1220	GLY	-	expression tag	UNP P0DTC2
D	1221	GLN	-	expression tag	UNP P0DTC2
D	1222	ALA	-	expression tag	UNP P0DTC2
D	1223	TYR	-	expression tag	UNP P0DTC2
D	1224	VAL	-	expression tag	UNP P0DTC2
D	1225	ARG	-	expression tag	UNP P0DTC2
D	1226	LYS	-	expression tag	UNP P0DTC2
D	1227	ASP	-	expression tag	UNP P0DTC2
D	1228	GLY	-	expression tag	UNP P0DTC2
D	1229	GLU	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	1230	TRP	-	expression tag	UNP P0DTC2
D	1231	VAL	-	expression tag	UNP P0DTC2
D	1232	LEU	-	expression tag	UNP P0DTC2
D	1233	LEU	-	expression tag	UNP P0DTC2
D	1234	SER	-	expression tag	UNP P0DTC2
D	1235	THR	-	expression tag	UNP P0DTC2
D	1236	PHE	-	expression tag	UNP P0DTC2
D	1237	LEU	-	expression tag	UNP P0DTC2
D	1238	GLU	-	expression tag	UNP P0DTC2
D	1239	ASN	-	expression tag	UNP P0DTC2
D	1240	LEU	-	expression tag	UNP P0DTC2
D	1241	TYR	-	expression tag	UNP P0DTC2
D	1242	PHE	-	expression tag	UNP P0DTC2
D	1243	GLN	-	expression tag	UNP P0DTC2
D	1244	GLY	-	expression tag	UNP P0DTC2
D	1245	ASP	-	expression tag	UNP P0DTC2
D	1246	TYR	-	expression tag	UNP P0DTC2
D	1247	LYS	-	expression tag	UNP P0DTC2
D	1248	ASP	-	expression tag	UNP P0DTC2
D	1249	ASP	-	expression tag	UNP P0DTC2
D	1250	ASP	-	expression tag	UNP P0DTC2
D	1251	ASP	-	expression tag	UNP P0DTC2
D	1252	LYS	-	expression tag	UNP P0DTC2
D	1253	HIS	-	expression tag	UNP P0DTC2
D	1254	HIS	-	expression tag	UNP P0DTC2
D	1255	HIS	-	expression tag	UNP P0DTC2
D	1256	HIS	-	expression tag	UNP P0DTC2
D	1257	HIS	-	expression tag	UNP P0DTC2
D	1258	HIS	-	expression tag	UNP P0DTC2
D	1259	HIS	-	expression tag	UNP P0DTC2
D	1260	HIS	-	expression tag	UNP P0DTC2
D	1261	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	142	ASP	GLY	variant	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	210A	ILE	-	insertion	UNP P0DTC2
C	210B	VAL	-	insertion	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	210C	ARG	ASN	conflict	UNP P0DTC2
C	210D	GLU	LEU	conflict	UNP P0DTC2
C	210E	PRO	VAL	conflict	UNP P0DTC2
C	210F	GLU	ARG	conflict	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
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	variant	UNP P0DTC2
C	683	SER	ARG	variant	UNP P0DTC2
C	685	SER	ARG	variant	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2
C	856	LYS	ASN	variant	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	variant	UNP P0DTC2
C	987	PRO	VAL	variant	UNP P0DTC2
C	1209	GLY	-	expression tag	UNP P0DTC2
C	1210	SER	-	expression tag	UNP P0DTC2
C	1211	GLY	-	expression tag	UNP P0DTC2
C	1212	TYR	-	expression tag	UNP P0DTC2
C	1213	ILE	-	expression tag	UNP P0DTC2
C	1214	PRO	-	expression tag	UNP P0DTC2
C	1215	GLU	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1216	ALA	-	expression tag	UNP P0DTC2
C	1217	PRO	-	expression tag	UNP P0DTC2
C	1218	ARG	-	expression tag	UNP P0DTC2
C	1219	ASP	-	expression tag	UNP P0DTC2
C	1220	GLY	-	expression tag	UNP P0DTC2
C	1221	GLN	-	expression tag	UNP P0DTC2
C	1222	ALA	-	expression tag	UNP P0DTC2
C	1223	TYR	-	expression tag	UNP P0DTC2
C	1224	VAL	-	expression tag	UNP P0DTC2
C	1225	ARG	-	expression tag	UNP P0DTC2
C	1226	LYS	-	expression tag	UNP P0DTC2
C	1227	ASP	-	expression tag	UNP P0DTC2
C	1228	GLY	-	expression tag	UNP P0DTC2
C	1229	GLU	-	expression tag	UNP P0DTC2
C	1230	TRP	-	expression tag	UNP P0DTC2
C	1231	VAL	-	expression tag	UNP P0DTC2
C	1232	LEU	-	expression tag	UNP P0DTC2
C	1233	LEU	-	expression tag	UNP P0DTC2
C	1234	SER	-	expression tag	UNP P0DTC2
C	1235	THR	-	expression tag	UNP P0DTC2
C	1236	PHE	-	expression tag	UNP P0DTC2
C	1237	LEU	-	expression tag	UNP P0DTC2
C	1238	GLU	-	expression tag	UNP P0DTC2
C	1239	ASN	-	expression tag	UNP P0DTC2
C	1240	LEU	-	expression tag	UNP P0DTC2
C	1241	TYR	-	expression tag	UNP P0DTC2
C	1242	PHE	-	expression tag	UNP P0DTC2
C	1243	GLN	-	expression tag	UNP P0DTC2
C	1244	GLY	-	expression tag	UNP P0DTC2
C	1245	ASP	-	expression tag	UNP P0DTC2
C	1246	TYR	-	expression tag	UNP P0DTC2
C	1247	LYS	-	expression tag	UNP P0DTC2
C	1248	ASP	-	expression tag	UNP P0DTC2
C	1249	ASP	-	expression tag	UNP P0DTC2
C	1250	ASP	-	expression tag	UNP P0DTC2
C	1251	ASP	-	expression tag	UNP P0DTC2
C	1252	LYS	-	expression tag	UNP P0DTC2
C	1253	HIS	-	expression tag	UNP P0DTC2
C	1254	HIS	-	expression tag	UNP P0DTC2
C	1255	HIS	-	expression tag	UNP P0DTC2
C	1256	HIS	-	expression tag	UNP P0DTC2
C	1257	HIS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1258	HIS	-	expression tag	UNP P0DTC2
C	1259	HIS	-	expression tag	UNP P0DTC2
C	1260	HIS	-	expression tag	UNP P0DTC2
C	1261	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	142	ASP	GLY	variant	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	210A	ILE	-	insertion	UNP P0DTC2
B	210B	VAL	-	insertion	UNP P0DTC2
B	210C	ARG	ASN	conflict	UNP P0DTC2
B	210D	GLU	LEU	conflict	UNP P0DTC2
B	210E	PRO	VAL	conflict	UNP P0DTC2
B	210F	GLU	ARG	conflict	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
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	variant	UNP P0DTC2
B	683	SER	ARG	variant	UNP P0DTC2
B	685	SER	ARG	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	856	LYS	ASN	variant	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	variant	UNP P0DTC2
B	987	PRO	VAL	variant	UNP P0DTC2
B	1209	GLY	-	expression tag	UNP P0DTC2
B	1210	SER	-	expression tag	UNP P0DTC2
B	1211	GLY	-	expression tag	UNP P0DTC2
B	1212	TYR	-	expression tag	UNP P0DTC2
B	1213	ILE	-	expression tag	UNP P0DTC2
B	1214	PRO	-	expression tag	UNP P0DTC2
B	1215	GLU	-	expression tag	UNP P0DTC2
B	1216	ALA	-	expression tag	UNP P0DTC2
B	1217	PRO	-	expression tag	UNP P0DTC2
B	1218	ARG	-	expression tag	UNP P0DTC2
B	1219	ASP	-	expression tag	UNP P0DTC2
B	1220	GLY	-	expression tag	UNP P0DTC2
B	1221	GLN	-	expression tag	UNP P0DTC2
B	1222	ALA	-	expression tag	UNP P0DTC2
B	1223	TYR	-	expression tag	UNP P0DTC2
B	1224	VAL	-	expression tag	UNP P0DTC2
B	1225	ARG	-	expression tag	UNP P0DTC2
B	1226	LYS	-	expression tag	UNP P0DTC2
B	1227	ASP	-	expression tag	UNP P0DTC2
B	1228	GLY	-	expression tag	UNP P0DTC2
B	1229	GLU	-	expression tag	UNP P0DTC2
B	1230	TRP	-	expression tag	UNP P0DTC2
B	1231	VAL	-	expression tag	UNP P0DTC2
B	1232	LEU	-	expression tag	UNP P0DTC2
B	1233	LEU	-	expression tag	UNP P0DTC2
B	1234	SER	-	expression tag	UNP P0DTC2
B	1235	THR	-	expression tag	UNP P0DTC2
B	1236	PHE	-	expression tag	UNP P0DTC2
B	1237	LEU	-	expression tag	UNP P0DTC2
B	1238	GLU	-	expression tag	UNP P0DTC2
B	1239	ASN	-	expression tag	UNP P0DTC2
B	1240	LEU	-	expression tag	UNP P0DTC2
B	1241	TYR	-	expression tag	UNP P0DTC2
B	1242	PHE	-	expression tag	UNP P0DTC2
B	1243	GLN	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1244	GLY	-	expression tag	UNP P0DTC2
B	1245	ASP	-	expression tag	UNP P0DTC2
B	1246	TYR	-	expression tag	UNP P0DTC2
B	1247	LYS	-	expression tag	UNP P0DTC2
B	1248	ASP	-	expression tag	UNP P0DTC2
B	1249	ASP	-	expression tag	UNP P0DTC2
B	1250	ASP	-	expression tag	UNP P0DTC2
B	1251	ASP	-	expression tag	UNP P0DTC2
B	1252	LYS	-	expression tag	UNP P0DTC2
B	1253	HIS	-	expression tag	UNP P0DTC2
B	1254	HIS	-	expression tag	UNP P0DTC2
B	1255	HIS	-	expression tag	UNP P0DTC2
B	1256	HIS	-	expression tag	UNP P0DTC2
B	1257	HIS	-	expression tag	UNP P0DTC2
B	1258	HIS	-	expression tag	UNP P0DTC2
B	1259	HIS	-	expression tag	UNP P0DTC2
B	1260	HIS	-	expression tag	UNP P0DTC2
B	1261	HIS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called Angiotensin-converting enzyme 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	597	Total	C	N	O	S	0	0
			4870	3115	806	920	29		

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q9BYF1
A	1	HIS	-	expression tag	UNP Q9BYF1
A	2	SER	-	expression tag	UNP Q9BYF1
A	3	SER	-	expression tag	UNP Q9BYF1
A	4	ALA	-	expression tag	UNP Q9BYF1
A	5	LEU	-	expression tag	UNP Q9BYF1
A	6	LEU	-	expression tag	UNP Q9BYF1
A	7	CYS	-	expression tag	UNP Q9BYF1
A	8	CYS	-	expression tag	UNP Q9BYF1
A	9	LEU	-	expression tag	UNP Q9BYF1
A	10	VAL	-	expression tag	UNP Q9BYF1
A	11	LEU	-	expression tag	UNP Q9BYF1
A	12	LEU	-	expression tag	UNP Q9BYF1
A	13	THR	-	expression tag	UNP Q9BYF1

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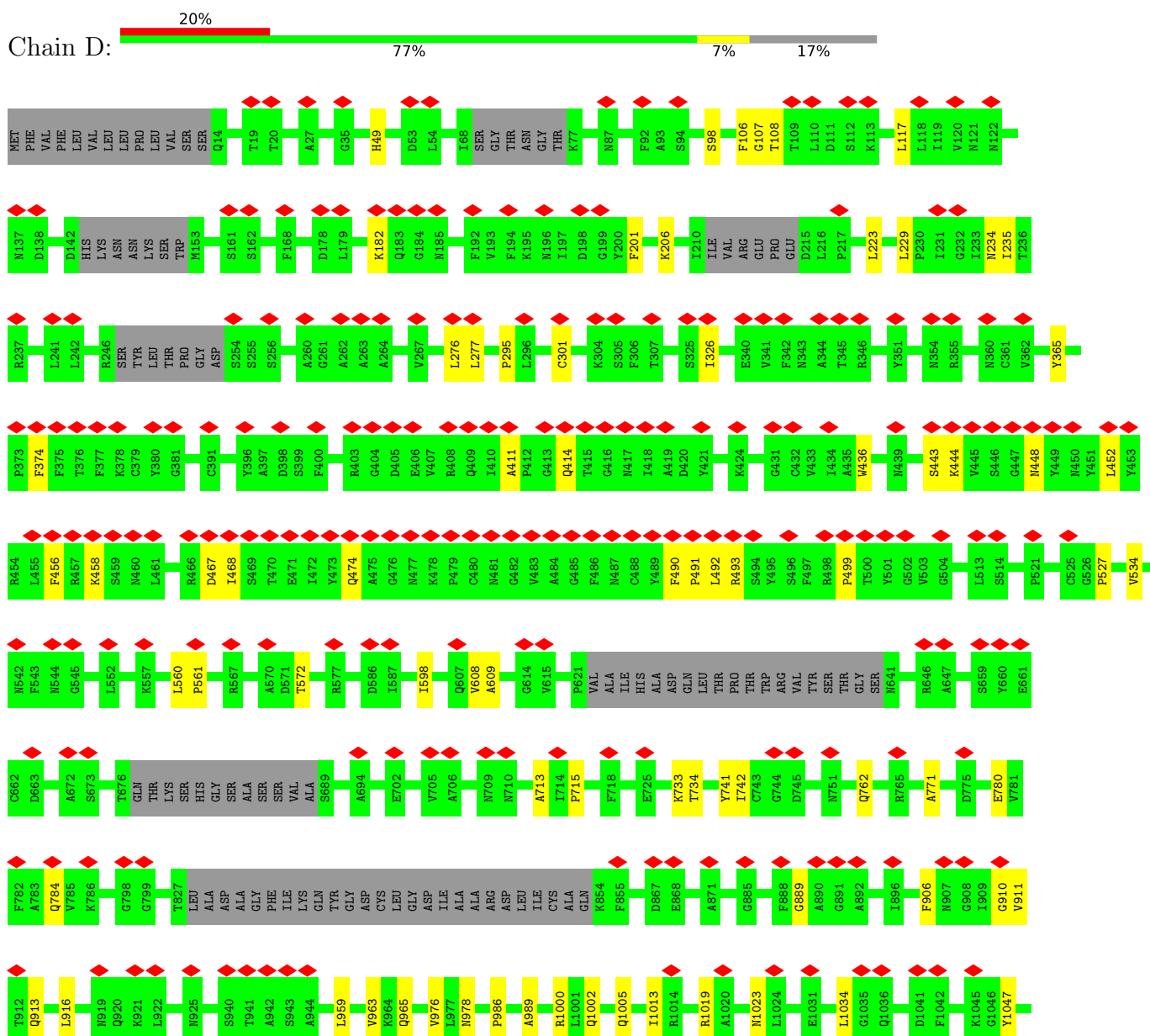
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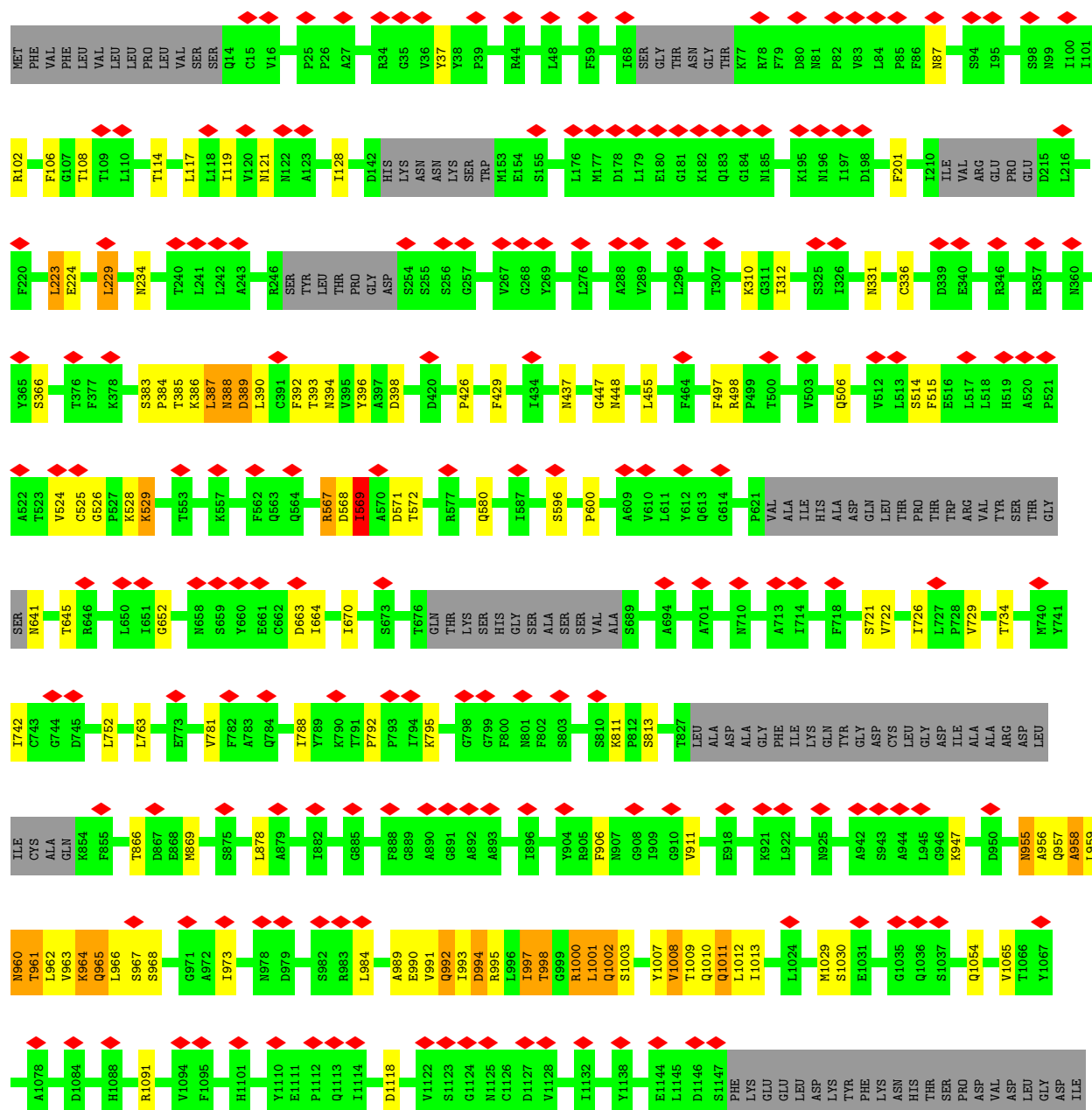
Chain	Residue	Modelled	Actual	Comment	Reference
A	14	GLY	-	expression tag	UNP Q9BYF1
A	15	VAL	-	expression tag	UNP Q9BYF1
A	16	ARG	-	expression tag	UNP Q9BYF1
A	616	HIS	-	expression tag	UNP Q9BYF1
A	617	HIS	-	expression tag	UNP Q9BYF1
A	618	HIS	-	expression tag	UNP Q9BYF1
A	619	HIS	-	expression tag	UNP Q9BYF1
A	620	HIS	-	expression tag	UNP Q9BYF1
A	621	HIS	-	expression tag	UNP Q9BYF1
A	622	HIS	-	expression tag	UNP Q9BYF1
A	623	HIS	-	expression tag	UNP Q9BYF1
A	624	HIS	-	expression tag	UNP Q9BYF1

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

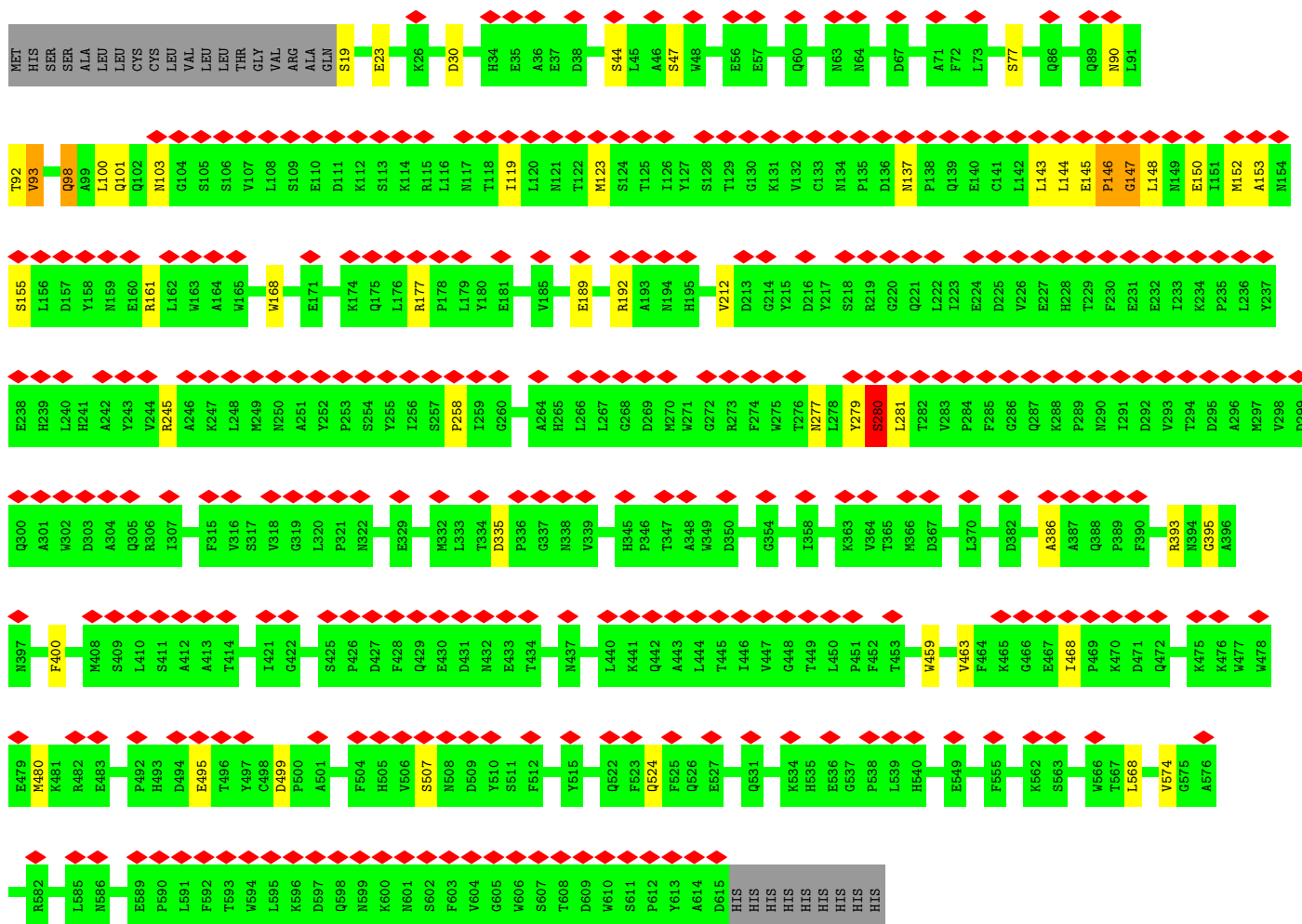
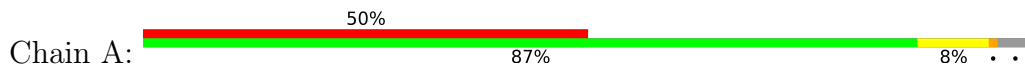
• Molecule 1: Spike glycoprotein







- Molecule 2: Angiotensin-converting enzyme 2



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	21125	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.981	Depositor
Minimum map value	-0.000	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.034	Depositor
Recommended contour level	0.12	Depositor
Map size (Å)	393.48, 393.48, 393.48	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.093, 1.093, 1.093	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	B	0.56	19/8410 (0.2%)	0.78	26/11435 (0.2%)
1	C	0.38	6/8410 (0.1%)	0.62	7/11435 (0.1%)
1	D	0.29	0/8410	0.57	1/11435 (0.0%)
2	A	0.40	4/5007 (0.1%)	0.68	4/6803 (0.1%)
All	All	0.42	29/30237 (0.1%)	0.66	38/41108 (0.1%)

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	335	ASP	C-N	7.80	1.52	1.33
1	B	1003	SER	CA-C	-7.38	1.43	1.52
2	A	499	ASP	C-N	7.28	1.42	1.34
1	B	336	CYS	C-N	7.23	1.50	1.33
1	B	997	ILE	CA-C	-7.03	1.43	1.52
1	B	960	ASN	CA-C	-6.82	1.44	1.52
1	B	1001	LEU	CA-C	-6.70	1.43	1.52
1	B	955	ASN	C-O	6.41	1.31	1.24
1	C	25	PRO	C-N	6.32	1.48	1.33
1	B	1007	TYR	CA-C	-6.01	1.45	1.52
1	B	995	ARG	CA-C	-6.00	1.45	1.52
1	B	229	LEU	C-N	5.97	1.47	1.33
1	C	994	ASP	CA-C	-5.87	1.45	1.52
1	C	1004	LEU	CA-C	-5.83	1.45	1.52
1	B	956	ALA	CA-C	-5.70	1.44	1.52
1	C	1003	SER	CA-C	-5.65	1.45	1.52
1	B	1000	ARG	CA-C	-5.64	1.45	1.52
2	A	92	THR	C-O	5.56	1.30	1.24
1	B	964	LYS	CA-C	-5.47	1.45	1.52
1	B	957	GLN	CA-C	-5.42	1.45	1.52
1	B	958	ALA	CA-C	-5.39	1.46	1.52
2	A	277	ASN	N-CA	5.39	1.53	1.46
1	C	1005	GLN	N-CA	-5.20	1.40	1.46
1	B	998	THR	CA-C	-5.17	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	994	ASP	CA-C	-5.17	1.46	1.52
1	B	1002	GLN	N-CA	-5.15	1.40	1.46
1	B	961	THR	N-CA	-5.12	1.40	1.46
1	C	995	ARG	N-CA	-5.10	1.40	1.46
1	B	1008	VAL	N-CA	-5.05	1.40	1.46

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	526	GLY	CA-C-O	12.69	139.66	121.52
1	B	957	GLN	N-CA-C	-12.15	98.04	111.28
1	B	526	GLY	N-CA-C	10.49	133.74	112.34
1	B	389	ASP	N-CA-C	9.58	121.32	111.07
1	C	1002	GLN	N-CA-C	-9.47	100.82	112.38
1	B	526	GLY	O-C-N	-9.41	112.36	121.77
1	B	569	ILE	N-CA-C	-9.14	101.30	110.62
1	B	390	LEU	N-CA-C	8.41	121.06	109.11
1	C	620	VAL	N-CA-C	7.99	121.94	112.35
1	B	526	GLY	CA-C-N	-7.58	111.56	120.13
1	B	526	GLY	C-N-CA	-7.58	111.56	120.13
1	B	524	VAL	N-CA-C	7.43	118.53	107.15
1	C	618	THR	N-CA-C	-7.43	103.27	111.36
1	C	223	LEU	N-CA-C	7.39	120.30	108.41
1	B	1002	GLN	N-CA-C	-6.96	103.69	111.28
1	B	528	LYS	CA-C-O	-6.61	113.36	121.89
1	B	525	CYS	N-CA-C	6.49	120.13	109.94
1	B	955	ASN	CA-C-O	6.32	127.26	120.24
2	A	101	GLN	N-CA-C	6.28	118.20	111.36
1	C	993	ILE	N-CA-C	6.17	116.95	110.72
2	A	98	GLN	N-CA-C	-6.05	104.69	111.28
1	B	1001	LEU	N-CA-C	-5.91	104.96	111.82
1	B	387	LEU	N-CA-C	-5.83	104.92	111.28
1	B	567	ARG	N-CA-C	5.79	118.03	109.69
1	B	385	THR	N-CA-C	-5.54	105.32	111.36
1	C	994	ASP	N-CA-C	-5.50	105.28	111.28
1	C	468	ILE	N-CA-C	-5.44	107.44	112.83
1	B	1007	TYR	CA-C-N	-5.37	113.10	120.46
1	B	1007	TYR	C-N-CA	-5.37	113.10	120.46
1	B	389	ASP	CA-C-N	-5.36	115.61	122.79
1	B	389	ASP	C-N-CA	-5.36	115.61	122.79
2	A	280	SER	CA-C-N	5.36	131.78	121.54
2	A	280	SER	C-N-CA	5.36	131.78	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	468	ILE	N-CA-C	-5.30	107.58	112.83
1	B	992	GLN	N-CA-C	-5.23	105.75	111.82
1	B	224	GLU	CA-C-N	-5.06	114.57	119.78
1	B	224	GLU	C-N-CA	-5.06	114.57	119.78
1	B	388	ASN	N-CA-C	-5.06	105.77	111.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	8219	0	8052	83	0
1	C	8219	0	8052	65	0
1	D	8219	0	8052	46	0
2	A	4870	0	4643	34	0
All	All	29527	0	28799	219	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 4.

All (219) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:TYR:HB3	1:C:223:LEU:HD21	1.54	0.88
1:C:990:GLU:HA	1:C:993:ILE:HG12	1.68	0.76
1:C:996:LEU:HD12	1:C:997:ILE:HG23	1.68	0.75
1:B:966:LEU:HG	1:B:1000:ARG:HH12	1.51	0.74
2:A:280:SER:O	2:A:281:LEU:HD12	1.88	0.72
1:C:993:ILE:HA	1:C:996:LEU:HG	1.72	0.71
1:B:387:LEU:HD22	1:B:515:PHE:HE2	1.60	0.67
1:B:734:THR:HG21	1:B:959:LEU:HD11	1.78	0.66
1:B:763:LEU:HD22	1:B:1008:VAL:HG11	1.79	0.65
1:C:1001:LEU:O	1:C:1002:GLN:C	2.39	0.64
2:A:155:SER:OG	2:A:161:ARG:HB3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1001:LEU:C	1:C:1003:SER:N	2.48	0.61
2:A:144:LEU:HA	2:A:148:LEU:HD22	1.84	0.59
1:B:964:LYS:C	1:B:966:LEU:H	2.11	0.59
1:B:966:LEU:C	1:B:968:SER:H	2.11	0.59
1:B:448:ASN:HB3	1:B:497:PHE:CD1	2.39	0.57
1:B:569:ILE:C	1:B:571:ASP:H	2.10	0.57
1:B:392:PHE:CG	1:B:515:PHE:HB3	2.39	0.57
1:B:1009:THR:HA	1:B:1012:LEU:HG	1.87	0.57
1:D:572:THR:HG22	1:C:856:LYS:HE2	1.87	0.56
1:B:386:LYS:HD2	1:B:389:ASP:HB2	1.87	0.56
1:B:964:LYS:O	1:B:966:LEU:N	2.39	0.56
1:B:958:ALA:O	1:B:961:THR:HB	2.06	0.56
2:A:123:MET:HB3	2:A:507:SER:HB3	1.88	0.56
1:D:106:PHE:HB2	1:D:117:LEU:HB3	1.86	0.56
1:B:1001:LEU:O	1:B:1002:GLN:C	2.48	0.56
1:D:201:PHE:HB2	1:D:229:LEU:HB2	1.87	0.55
1:C:617:CYS:O	1:C:620:VAL:HG22	2.06	0.55
1:D:762:GLN:HE21	1:B:961:THR:HG23	1.71	0.55
1:B:312:ILE:HD11	1:B:596:SER:HB3	1.89	0.55
1:B:1010:GLN:O	1:B:1011:GLN:C	2.47	0.55
1:C:106:PHE:HB3	1:C:235:ILE:HG12	1.88	0.55
2:A:279:TYR:C	2:A:281:LEU:H	2.14	0.55
1:C:1005:GLN:C	1:C:1007:TYR:N	2.65	0.54
1:B:331:ASN:HB2	1:B:580:GLN:HG3	1.88	0.54
1:B:962:LEU:O	1:B:965:GLN:HB3	2.06	0.54
1:C:201:PHE:HB2	1:C:229:LEU:HB2	1.89	0.54
1:C:1005:GLN:HA	1:C:1008:VAL:HG12	1.89	0.54
1:B:384:PRO:HA	1:B:387:LEU:HD12	1.89	0.54
1:C:189:LEU:HD13	1:C:217:PRO:HG2	1.89	0.53
1:B:993:ILE:O	1:B:997:ILE:HG12	2.08	0.53
1:D:326:ILE:HD11	1:D:534:VAL:HG12	1.91	0.53
1:D:1052:PHE:HB2	1:D:1063:LEU:HB2	1.90	0.52
1:C:102:ARG:NH1	1:C:177:MET:SD	2.82	0.52
2:A:143:LEU:HG	2:A:144:LEU:H	1.74	0.52
1:C:106:PHE:HB2	1:C:117:LEU:HB3	1.91	0.52
1:B:201:PHE:HB2	1:B:229:LEU:HB2	1.92	0.52
1:B:366:SER:HB2	1:B:388:ASN:HD21	1.74	0.52
1:C:21:ARG:HG3	1:C:79:PHE:HB2	1.92	0.52
1:B:119:ILE:HG12	1:B:128:ILE:HG12	1.91	0.52
2:A:152:MET:O	2:A:153:ALA:C	2.53	0.52
1:B:964:LYS:C	1:B:966:LEU:N	2.62	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:966:LEU:C	1:B:968:SER:N	2.67	0.51
1:C:443:SER:HB2	1:C:499:PRO:HD3	1.92	0.51
1:D:598:ILE:HB	1:D:609:ALA:HB3	1.92	0.51
1:C:669:GLY:HA2	1:C:697:MET:HE2	1.93	0.51
1:C:734:THR:HG21	1:C:959:LEU:HD11	1.92	0.51
1:B:878:LEU:HD21	1:B:1054:GLN:HE22	1.76	0.51
1:D:1013:ILE:HD13	1:C:1012:LEU:HD13	1.92	0.51
1:C:792:PRO:O	1:C:795:LYS:NZ	2.44	0.51
1:C:1005:GLN:O	1:C:1006:THR:C	2.52	0.51
1:B:792:PRO:O	1:B:795:LYS:NZ	2.43	0.50
2:A:144:LEU:HB2	2:A:168:TRP:CH2	2.46	0.50
1:B:1091:ARG:NH2	1:B:1118:ASP:O	2.46	0.49
1:B:955:ASN:O	1:B:958:ALA:HB3	2.11	0.49
1:C:990:GLU:HA	1:C:993:ILE:CG1	2.42	0.49
1:C:1019:ARG:NH1	1:C:1023:ASN:OD1	2.45	0.49
1:B:726:ILE:O	1:B:947:LYS:NZ	2.46	0.49
1:D:976:VAL:HG12	1:D:978:ASN:H	1.77	0.49
1:C:452:LEU:HD23	1:C:492:LEU:HB3	1.94	0.49
2:A:459:TRP:HB2	2:A:480:MET:HE1	1.95	0.49
2:A:144:LEU:HA	2:A:148:LEU:CD2	2.42	0.48
1:B:811:LYS:NZ	1:B:813:SER:OG	2.47	0.48
2:A:145:GLU:N	2:A:146:PRO:HD2	2.28	0.48
1:D:742:ILE:O	1:D:1000:ARG:NH1	2.47	0.48
1:C:98:SER:HB3	1:C:182:LYS:H	1.77	0.48
1:C:617:CYS:HA	1:C:620:VAL:HG22	1.94	0.48
2:A:144:LEU:HD13	2:A:168:TRP:CZ2	2.49	0.48
1:D:713:ALA:HB3	1:C:895:GLN:H	1.78	0.48
1:D:1019:ARG:NH1	1:D:1023:ASN:OD1	2.47	0.48
1:B:102:ARG:HD3	1:B:121:ASN:H	1.78	0.47
2:A:212:VAL:HG21	2:A:568:LEU:HD11	1.94	0.47
1:C:426:PRO:HG2	1:C:429:PHE:HB2	1.96	0.47
2:A:143:LEU:HB3	2:A:146:PRO:CD	2.44	0.47
2:A:147:GLY:O	2:A:148:LEU:C	2.57	0.47
1:D:444:LYS:HD3	1:D:448:ASN:HA	1.96	0.47
1:D:452:LEU:HD23	1:D:492:LEU:HB3	1.95	0.47
1:D:913:GLN:HA	1:D:916:LEU:HD13	1.97	0.47
1:C:126:VAL:HB	1:C:172:SER:HB2	1.97	0.47
1:C:328:ARG:NH1	1:C:578:ASP:OD1	2.48	0.47
1:B:645:THR:HG21	1:B:670:ILE:HD13	1.96	0.47
2:A:146:PRO:O	2:A:148:LEU:N	2.47	0.47
2:A:245:ARG:HD3	2:A:258:PRO:HA	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:326:ILE:HD11	1:C:534:VAL:HG12	1.96	0.47
1:C:411:ALA:HB3	1:C:414:GLN:HB2	1.96	0.47
1:D:443:SER:HB2	1:D:499:PRO:HD3	1.97	0.47
2:A:19:SER:N	2:A:23:GLU:OE1	2.49	0.46
1:D:374:PHE:HA	1:D:436:TRP:HD1	1.80	0.46
1:D:733:LYS:HE3	1:D:771:ALA:HB2	1.97	0.46
1:C:1001:LEU:C	1:C:1003:SER:H	2.20	0.46
1:D:741:TYR:OH	1:D:965:GLN:NE2	2.49	0.46
1:B:569:ILE:C	1:B:571:ASP:N	2.68	0.46
1:B:742:ILE:HA	1:B:1000:ARG:HE	1.81	0.46
1:D:107:GLY:H	1:D:235:ILE:HG23	1.81	0.46
1:D:411:ALA:HB3	1:D:414:GLN:HB2	1.97	0.46
1:C:107:GLY:H	1:C:235:ILE:HG23	1.81	0.46
1:B:781:VAL:HG13	1:B:1029:MET:HG3	1.98	0.46
1:B:1009:THR:HA	1:B:1012:LEU:CG	2.46	0.46
1:D:734:THR:HG21	1:D:959:LEU:HD21	1.97	0.46
1:B:966:LEU:O	1:B:968:SER:N	2.48	0.46
1:D:780:GLU:O	1:D:784:GLN:NE2	2.49	0.46
1:C:203:ILE:HB	1:C:227:VAL:HB	1.97	0.46
1:C:819:GLU:HG2	1:C:1055:SER:HB3	1.98	0.46
2:A:119:ILE:HG12	2:A:123:MET:HE3	1.98	0.46
1:D:910:GLY:O	1:D:1106:GLN:NE2	2.49	0.45
1:C:908:GLY:O	1:C:1038:LYS:NZ	2.49	0.45
1:B:310:LYS:HG3	1:B:600:PRO:HA	1.99	0.45
1:B:906:PHE:HB3	1:B:911:VAL:HB	1.99	0.45
2:A:145:GLU:N	2:A:146:PRO:CD	2.78	0.45
1:C:193:VAL:HB	1:C:204:TYR:HB2	1.99	0.45
1:D:98:SER:HB3	1:D:182:LYS:H	1.82	0.45
1:C:977:LEU:HD11	1:C:1000:ARG:HH12	1.82	0.45
1:B:398:ASP:OD1	1:B:398:ASP:N	2.50	0.45
1:D:490:PHE:O	1:D:493:ARG:NH1	2.49	0.45
1:D:906:PHE:HB3	1:D:911:VAL:HB	1.99	0.45
1:C:1004:LEU:H	1:C:1004:LEU:HG	1.53	0.45
1:B:991:VAL:O	1:B:992:GLN:C	2.55	0.45
1:B:990:GLU:HA	1:B:993:ILE:HG13	1.99	0.45
1:C:376:THR:HB	1:C:435:ALA:HB3	1.99	0.44
1:C:444:LYS:HD3	1:C:448:ASN:HA	1.98	0.44
1:C:598:ILE:HB	1:C:609:ALA:HB3	1.98	0.44
1:C:1004:LEU:HD12	1:C:1005:GLN:HG3	2.00	0.44
1:B:447:GLY:HA2	1:B:498:ARG:NH1	2.33	0.44
1:B:752:LEU:HD22	1:B:993:ILE:HG21	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:HIS:HB3	1:D:277:LEU:HD12	2.00	0.44
1:B:641:ASN:HB2	1:B:652:GLY:H	1.81	0.44
1:D:374:PHE:HD1	1:D:436:TRP:HB2	1.83	0.44
1:D:1089:PHE:HB3	1:C:913:GLN:HE21	1.81	0.44
1:C:762:GLN:HA	1:C:765:ARG:HG2	2.00	0.44
1:B:973:ILE:HD11	1:B:984:LEU:HD11	2.00	0.44
2:A:524:GLN:HG2	2:A:574:VAL:HG11	2.00	0.44
1:D:467:ASP:OD1	1:D:467:ASP:N	2.51	0.44
1:D:715:PRO:HA	1:D:1072:GLU:HA	1.99	0.44
1:B:965:GLN:HB3	1:B:965:GLN:HE21	1.55	0.44
1:C:994:ASP:O	1:C:995:ARG:C	2.61	0.43
1:B:310:LYS:HG2	1:B:664:ILE:HD11	2.00	0.43
1:B:383:SER:O	1:B:387:LEU:N	2.47	0.43
1:D:911:VAL:HA	1:D:1106:GLN:HE22	1.84	0.43
1:B:959:LEU:O	1:B:960:ASN:C	2.58	0.43
1:D:276:LEU:HD11	1:D:301:CYS:HA	2.01	0.43
1:C:1048:HIS:NE2	1:C:1051:SER:OG	2.42	0.43
1:C:910:GLY:HA3	1:C:1107:ARG:HH21	1.82	0.43
1:B:386:LYS:HA	1:B:389:ASP:HB2	2.00	0.43
1:B:437:ASN:ND2	1:B:506:GLN:OE1	2.48	0.43
2:A:137:ASN:OD1	2:A:137:ASN:N	2.52	0.43
1:B:866:THR:HG22	1:B:869:MET:HE3	1.99	0.43
2:A:386:ALA:O	2:A:393:ARG:NH2	2.52	0.43
2:A:395:GLY:HA3	2:A:400:PHE:HB3	2.00	0.43
1:C:560:LEU:HD12	1:C:561:PRO:HD2	2.01	0.43
1:B:396:TYR:HB2	1:B:514:SER:HB3	1.99	0.43
1:B:1010:GLN:O	1:B:1012:LEU:N	2.52	0.43
2:A:177:ARG:NE	2:A:495:GLU:O	2.51	0.43
1:C:456:PHE:HB2	1:C:491:PRO:HB3	2.01	0.43
1:C:1004:LEU:O	1:C:1008:VAL:N	2.47	0.43
1:C:364:ASP:OD1	1:C:364:ASP:N	2.51	0.42
1:C:1041:ASP:HB3	1:B:1030:SER:HB2	2.01	0.42
1:C:438:SER:HB3	1:C:442:ASP:HB2	2.00	0.42
2:A:146:PRO:O	2:A:147:GLY:C	2.62	0.42
1:C:365:TYR:H	1:C:527:PRO:HG3	1.83	0.42
1:C:994:ASP:OD1	1:C:994:ASP:N	2.53	0.42
1:D:456:PHE:HB2	1:D:491:PRO:HB3	2.00	0.42
1:B:663:ASP:OD1	1:B:663:ASP:N	2.52	0.42
1:C:490:PHE:O	1:C:493:ARG:NH1	2.49	0.42
1:B:106:PHE:HB2	1:B:117:LEU:HB3	2.00	0.42
1:B:108:THR:HB	1:B:114:THR:HG21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:722:VAL:HG22	1:B:1065:VAL:HG22	2.00	0.42
1:B:962:LEU:HA	1:B:962:LEU:HD12	1.78	0.42
1:B:721:SER:OG	1:B:722:VAL:N	2.53	0.41
1:D:458:LYS:NZ	1:D:474:GLN:O	2.53	0.41
1:C:950:ASP:O	1:C:954:HIS:ND1	2.45	0.41
2:A:524:GLN:HE21	2:A:574:VAL:HG21	1.84	0.41
1:C:699:LEU:HG	1:B:788:ILE:HG13	2.02	0.41
1:B:426:PRO:HG2	1:B:429:PHE:HB2	2.02	0.41
1:B:455:LEU:HD21	2:A:30:ASP:HB3	2.02	0.41
1:D:560:LEU:HD12	1:D:561:PRO:HD2	2.03	0.41
1:D:963:VAL:HG13	1:B:571:ASP:HB2	2.03	0.41
1:B:108:THR:OG1	1:B:234:ASN:O	2.38	0.41
1:D:889:GLY:HA3	1:D:1034:LEU:HD22	2.01	0.41
1:B:568:ASP:HB3	1:B:572:THR:H	1.85	0.41
2:A:90:ASN:HB3	2:A:93:VAL:HG12	2.02	0.41
1:D:206:LYS:HB2	1:D:223:LEU:HA	2.03	0.41
1:B:1009:THR:CA	1:B:1012:LEU:HG	2.50	0.41
2:A:77:SER:OG	2:A:103:ASN:OD1	2.39	0.41
1:C:80:ASP:OD1	1:C:80:ASP:N	2.52	0.41
1:B:497:PHE:O	1:B:498:ARG:HD3	2.20	0.41
2:A:143:LEU:HB3	2:A:146:PRO:HD2	2.03	0.41
1:D:986:PRO:HA	1:D:989:ALA:HB3	2.02	0.41
1:D:1002:GLN:HA	1:D:1005:GLN:HG2	2.02	0.41
1:C:1005:GLN:O	1:C:1007:TYR:N	2.53	0.41
1:B:37:TYR:CB	1:B:223:LEU:HB2	2.51	0.41
1:B:223:LEU:HD13	1:B:223:LEU:HA	1.67	0.41
1:B:392:PHE:CD2	1:B:515:PHE:HB3	2.56	0.41
1:B:567:ARG:HH21	1:B:571:ASP:C	2.29	0.41
1:B:1012:LEU:HD12	1:B:1013:ILE:HG13	2.02	0.41
2:A:44:SER:HA	2:A:47:SER:HB3	2.03	0.41
1:B:87:ASN:OD1	1:B:87:ASN:N	2.52	0.41
1:B:568:ASP:HB2	1:B:572:THR:O	2.20	0.41
2:A:463:VAL:HG12	2:A:468:ILE:HD11	2.01	0.41
1:B:392:PHE:C	1:B:394:ASN:N	2.80	0.40
1:D:365:TYR:H	1:D:527:PRO:HG3	1.86	0.40
1:C:1004:LEU:O	1:C:1007:TYR:HB3	2.20	0.40
1:D:108:THR:OG1	1:D:234:ASN:O	2.40	0.40
1:D:1047:TYR:HB2	1:D:1067:TYR:HB3	2.03	0.40
1:B:384:PRO:O	1:B:387:LEU:HB2	2.22	0.40
1:C:276:LEU:HD11	1:C:301:CYS:HA	2.04	0.40
1:B:994:ASP:O	1:B:998:THR:HG23	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:295:PRO:HD2	1:D:608:VAL:HG21	2.03	0.40
1:B:989:ALA:O	1:B:993:ILE:HG13	2.22	0.40
2:A:189:GLU:OE2	2:A:192:ARG:NH2	2.50	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	B	1032/1258 (82%)	967 (94%)	61 (6%)	4 (0%)	30	64
1	C	1032/1258 (82%)	981 (95%)	51 (5%)	0	100	100
1	D	1032/1258 (82%)	978 (95%)	54 (5%)	0	100	100
2	A	595/625 (95%)	566 (95%)	26 (4%)	3 (0%)	24	60
All	All	3691/4399 (84%)	3492 (95%)	192 (5%)	7 (0%)	44	75

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	529	LYS
2	A	280	SER
1	B	967	SER
1	B	1011	GLN
2	A	147	GLY
1	B	393	THR
2	A	146	PRO

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	B	918/1099 (84%)	912 (99%)	6 (1%)	76	79
1	C	918/1099 (84%)	914 (100%)	4 (0%)	84	83
1	D	918/1099 (84%)	918 (100%)	0	100	100
2	A	527/552 (96%)	522 (99%)	5 (1%)	70	76
All	All	3281/3849 (85%)	3266 (100%)	15 (0%)	78	82

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	994	ASP
1	C	996	LEU
1	C	1003	SER
1	C	1004	LEU
1	B	223	LEU
1	B	529	LYS
1	B	569	ILE
1	B	729	VAL
1	B	963	VAL
1	B	965	GLN
2	A	93	VAL
2	A	98	GLN
2	A	100	LEU
2	A	150	GLU
2	A	280	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	30	ASN
1	D	66	HIS
1	D	188	ASN
1	D	196	ASN
1	D	218	GLN
1	D	448	ASN
1	D	564	GLN
1	D	784	GLN

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Mol	Chain	Res	Type
1	D	960	ASN
1	D	1054	GLN
1	D	1088	HIS
1	D	1101	HIS
1	D	1106	GLN
1	D	1113	GLN
1	C	164	ASN
1	C	188	ASN
1	C	196	ASN
1	C	271	GLN
1	C	675	GLN
1	C	787	GLN
1	C	895	GLN
1	C	914	ASN
1	C	935	GLN
1	C	992	GLN
1	C	1011	GLN
1	C	1054	GLN
1	B	394	ASN
1	B	448	ASN
1	B	675	GLN
1	B	787	GLN
1	B	1005	GLN
1	B	1010	GLN
1	B	1054	GLN
1	B	1088	HIS
1	B	1101	HIS
2	A	24	GLN
2	A	53	ASN
2	A	76	GLN
2	A	96	GLN
2	A	277	ASN
2	A	330	ASN
2	A	417	HIS
2	A	472	GLN
2	A	508	ASN
2	A	524	GLN
2	A	526	GLN
2	A	535	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32857. These allow visual inspection of the internal detail of the map and identification of artifacts.

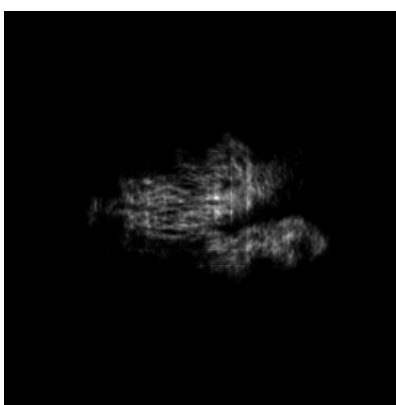
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

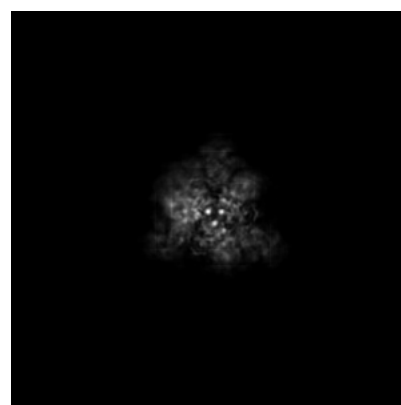
6.1.1 Primary map



X



Y



Z

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

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 180



Y Index: 180



Z Index: 180

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: 191



Y Index: 178

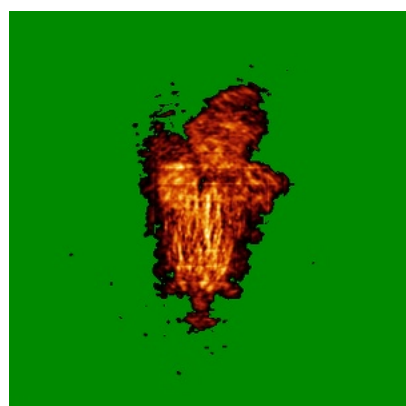


Z Index: 192

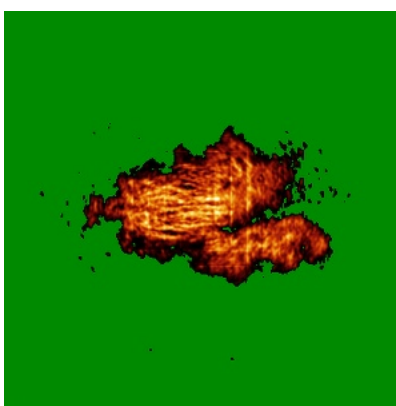
The images above show the largest variance slices of the map in three orthogonal directions.

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

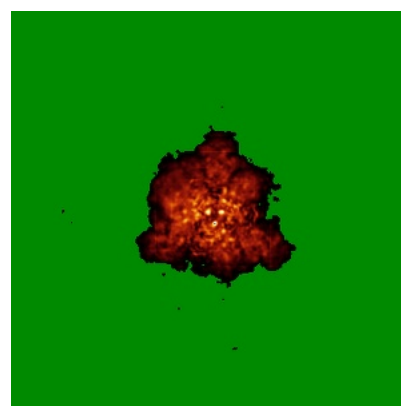
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.12. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

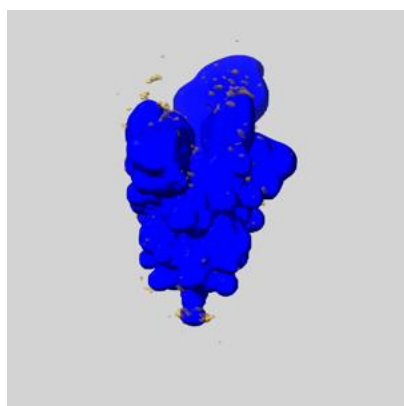
6.6 Mask visualisation [i](#)

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

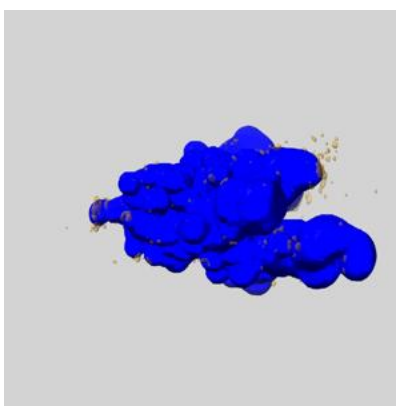
A mask typically either:

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

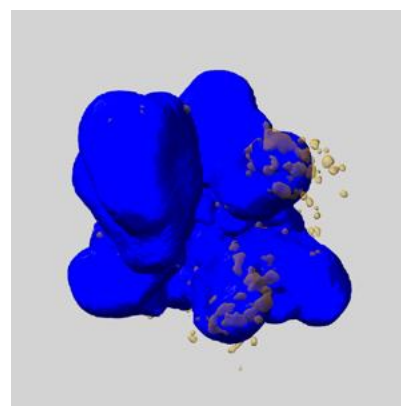
6.6.1 emd_32857_msk_1.map [i](#)



X



Y

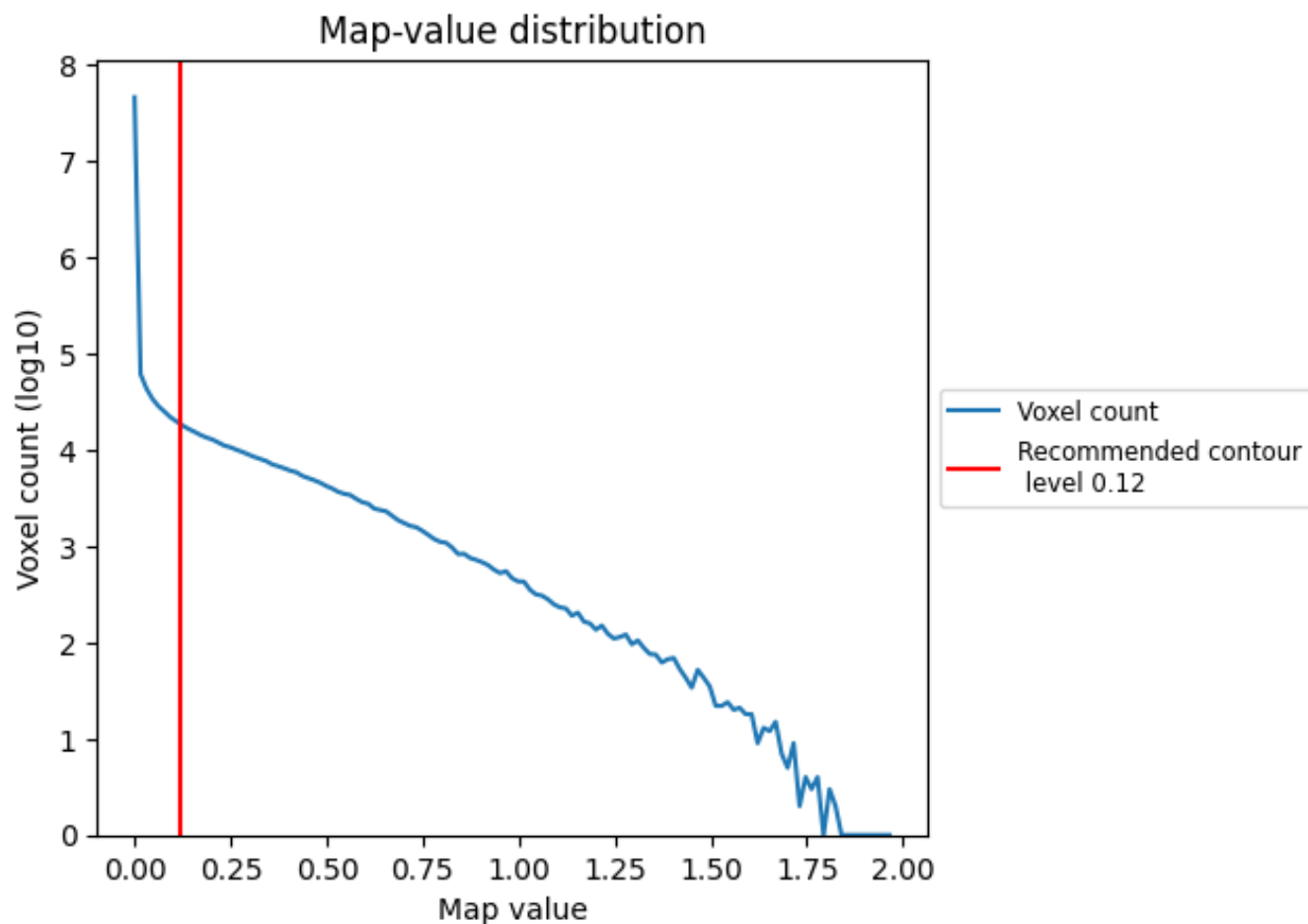


Z

7 Map analysis [i](#)

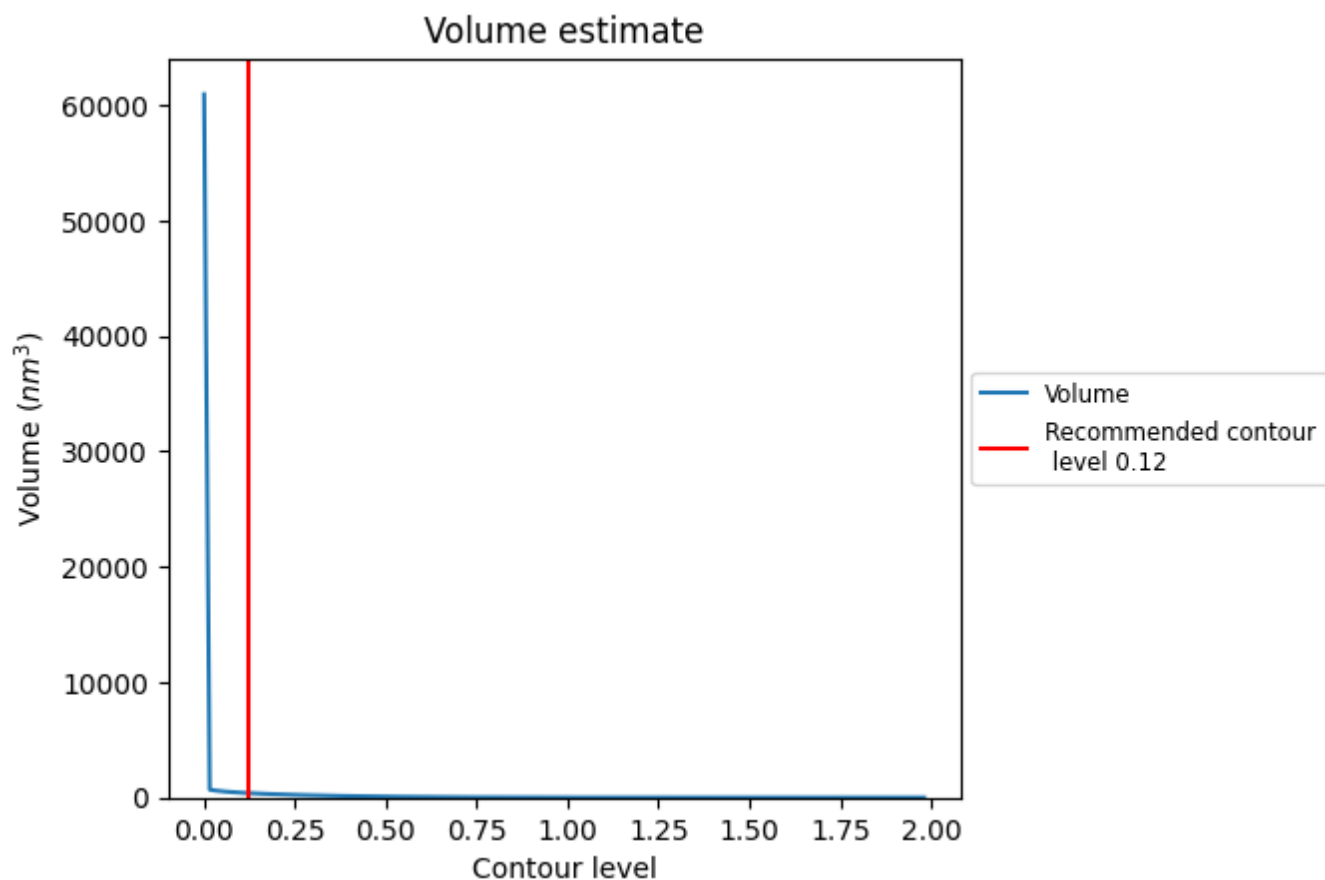
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

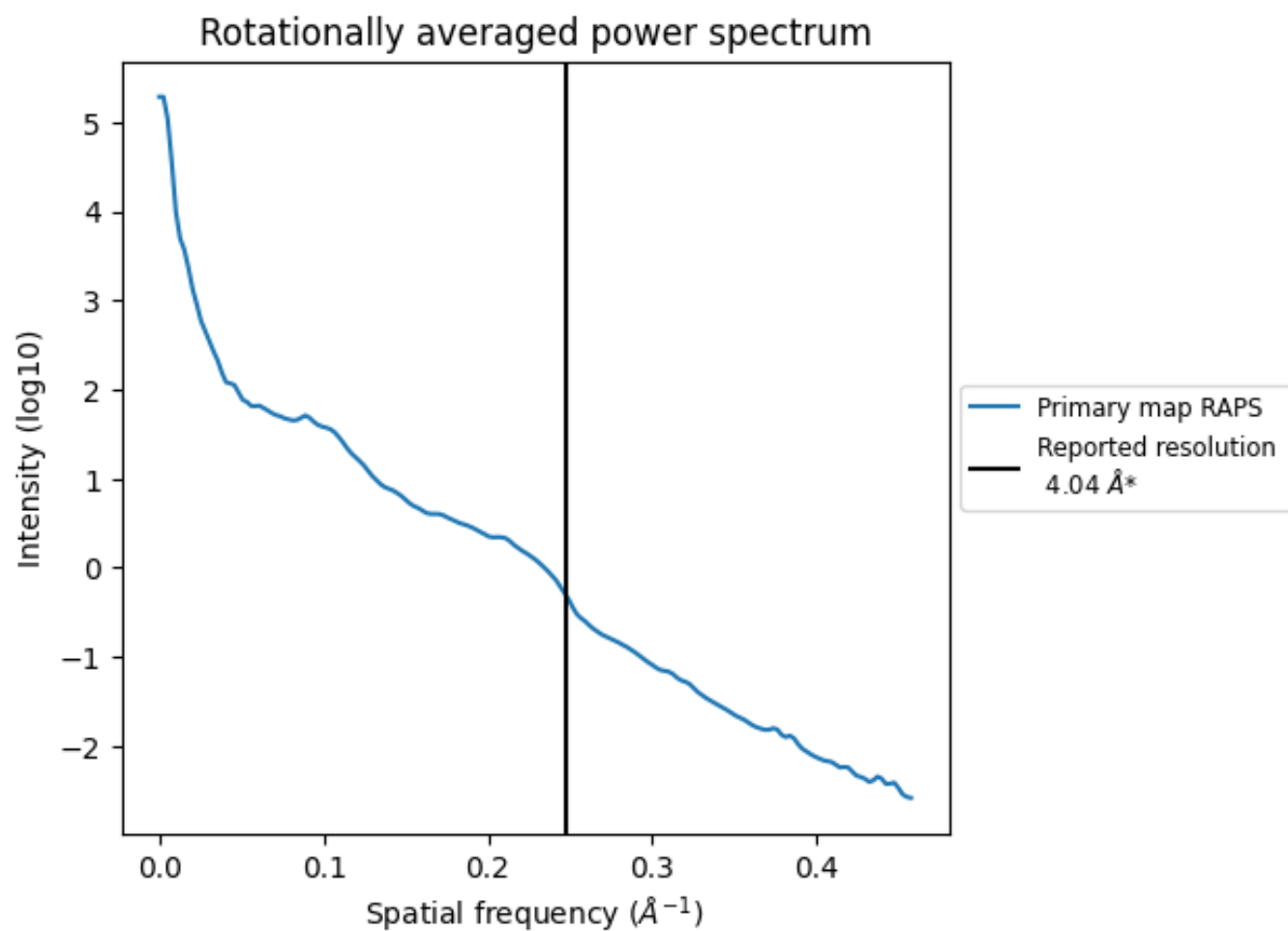
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 388 nm³; this corresponds to an approximate mass of 351 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 0.248 $\text{\AA}^{-1}$

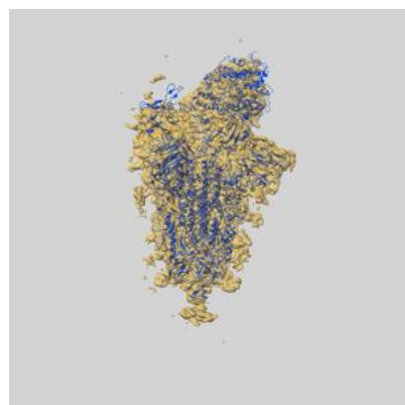
8 Fourier-Shell correlation ⓘ

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

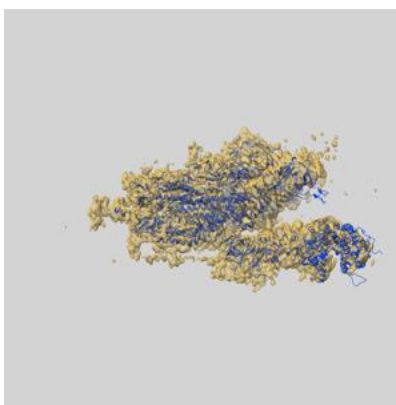
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-32857 and PDB model 7WVQ. Per-residue inclusion information can be found in section [3](#) on page [12](#).

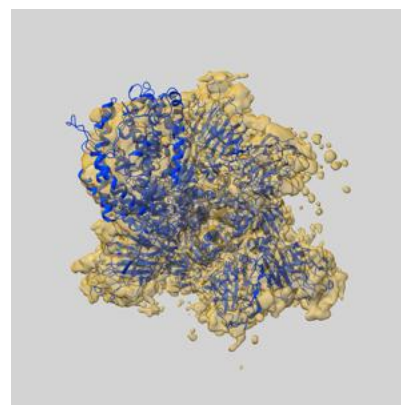
9.1 Map-model overlay [i](#)



X



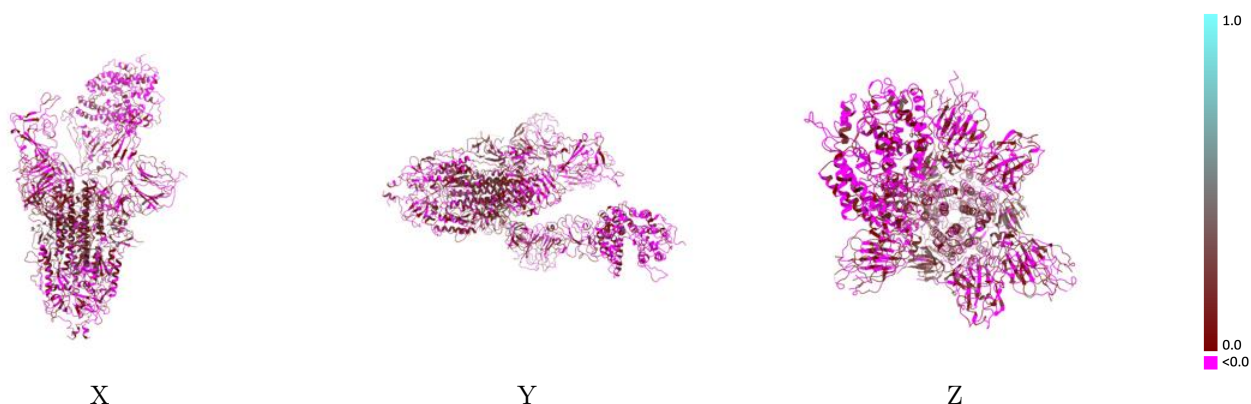
Y



Z

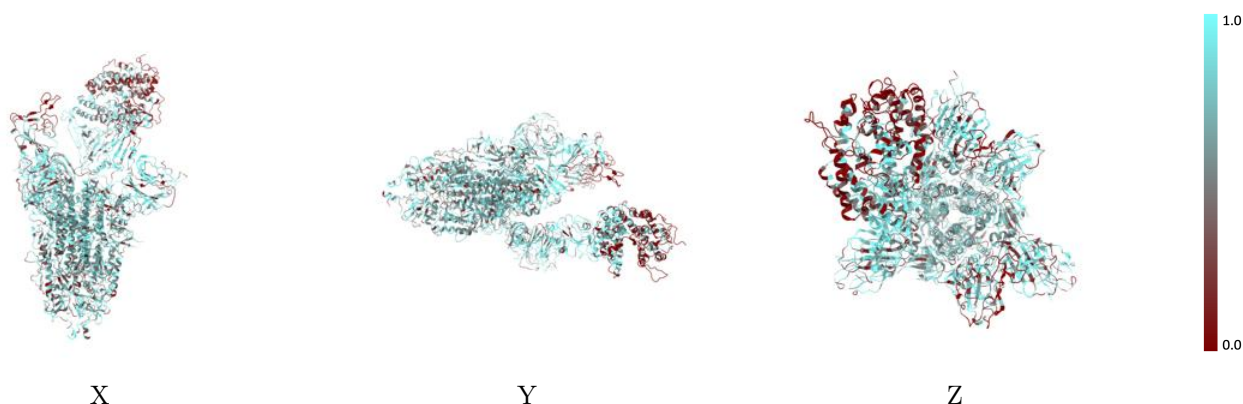
The images above show the 3D surface view of the map at the recommended contour level 0.12 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



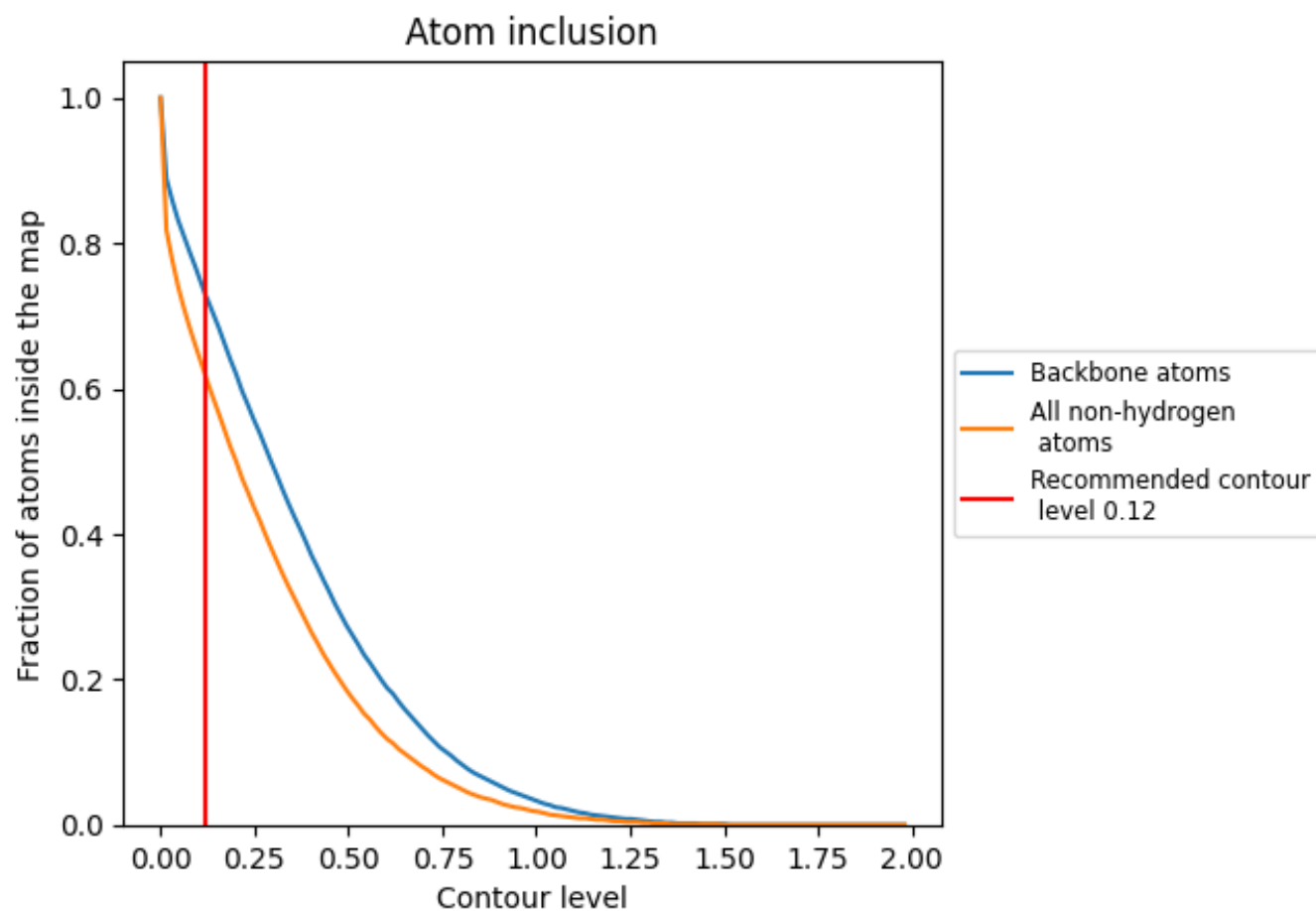
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.12).

9.4 Atom inclusion [i](#)



At the recommended contour level, 73% of all backbone atoms, 62% 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 (0.12) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6160	0.0570
A	0.4260	-0.0230
B	0.6880	0.0730
C	0.6370	0.0670
D	0.6340	0.0770

