



Full wwPDB EM Validation Report ⓘ

Mar 21, 2022 – 12:08 pm GMT

PDB ID : 7NY5
EMDB ID : EMD-12645
Title : Structure of SARS-CoV-2 Spike with P5C3 Fab
Authors : Perez, L.
Deposited on : 2021-03-20
Resolution : 3.70 Å(reported)
Based on initial model : 7K4N

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

EMDB validation analysis : 0.0.0.dev97
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.27

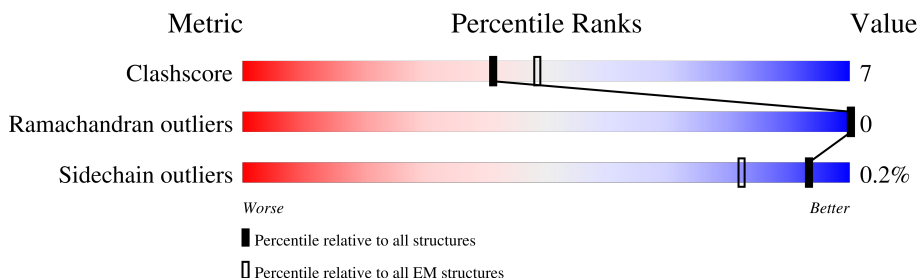
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.70 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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

Mol	Chain	Length	Quality of chain
1	A	1288	69% 9% 23%
1	E	1288	69% 9% 23%
1	G	1288	69% 9% 23%
2	B	215	21% 34% 16% 50%
2	D	215	23% 34% 16% 50%
2	H	215	22% 35% 15% 50%
3	C	231	8% 34% 19% 47%
3	F	231	7% 35% 18% 47%

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Mol	Chain	Length	Quality of chain
3	I	231	<div><div></div><div>8%</div><div>35%</div><div>18%</div><div>47%</div></div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 26466 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	998	Total	C	N	O	S	0	0
			7064	4529	1209	1292	34		
1	E	998	Total	C	N	O	S	0	0
			7064	4529	1209	1292	34		
1	G	998	Total	C	N	O	S	0	0
			7064	4529	1209	1292	34		

There are 267 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	682	GLY	ARG	conflict	UNP P0DTC2
A	683	SER	ARG	conflict	UNP P0DTC2
A	685	SER	ARG	conflict	UNP P0DTC2
A	817	PRO	PHE	conflict	UNP P0DTC2
A	892	PRO	ALA	conflict	UNP P0DTC2
A	899	PRO	ALA	conflict	UNP P0DTC2
A	942	PRO	ALA	conflict	UNP P0DTC2
A	986	PRO	LYS	conflict	UNP P0DTC2
A	987	PRO	VAL	conflict	UNP P0DTC2
A	1209	GLY	-	expression tag	UNP P0DTC2
A	1210	SER	-	expression tag	UNP P0DTC2
A	1211	GLY	-	expression tag	UNP P0DTC2
A	1212	TYR	-	expression tag	UNP P0DTC2
A	1213	ILE	-	expression tag	UNP P0DTC2
A	1214	PRO	-	expression tag	UNP P0DTC2
A	1215	GLU	-	expression tag	UNP P0DTC2
A	1216	ALA	-	expression tag	UNP P0DTC2
A	1217	PRO	-	expression tag	UNP P0DTC2
A	1218	ARG	-	expression tag	UNP P0DTC2
A	1219	ASP	-	expression tag	UNP P0DTC2
A	1220	GLY	-	expression tag	UNP P0DTC2
A	1221	GLN	-	expression tag	UNP P0DTC2
A	1222	ALA	-	expression tag	UNP P0DTC2
A	1223	TYR	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1224	VAL	-	expression tag	UNP P0DTC2
A	1225	ARG	-	expression tag	UNP P0DTC2
A	1226	LYS	-	expression tag	UNP P0DTC2
A	1227	ASP	-	expression tag	UNP P0DTC2
A	1228	GLY	-	expression tag	UNP P0DTC2
A	1229	GLU	-	expression tag	UNP P0DTC2
A	1230	TRP	-	expression tag	UNP P0DTC2
A	1231	VAL	-	expression tag	UNP P0DTC2
A	1232	LEU	-	expression tag	UNP P0DTC2
A	1233	LEU	-	expression tag	UNP P0DTC2
A	1234	SER	-	expression tag	UNP P0DTC2
A	1235	THR	-	expression tag	UNP P0DTC2
A	1236	PHE	-	expression tag	UNP P0DTC2
A	1237	LEU	-	expression tag	UNP P0DTC2
A	1238	GLY	-	expression tag	UNP P0DTC2
A	1239	ARG	-	expression tag	UNP P0DTC2
A	1240	SER	-	expression tag	UNP P0DTC2
A	1241	LEU	-	expression tag	UNP P0DTC2
A	1242	GLU	-	expression tag	UNP P0DTC2
A	1243	VAL	-	expression tag	UNP P0DTC2
A	1244	LEU	-	expression tag	UNP P0DTC2
A	1245	PHE	-	expression tag	UNP P0DTC2
A	1246	GLN	-	expression tag	UNP P0DTC2
A	1247	GLY	-	expression tag	UNP P0DTC2
A	1248	PRO	-	expression tag	UNP P0DTC2
A	1249	GLY	-	expression tag	UNP P0DTC2
A	1250	HIS	-	expression tag	UNP P0DTC2
A	1251	HIS	-	expression tag	UNP P0DTC2
A	1252	HIS	-	expression tag	UNP P0DTC2
A	1253	HIS	-	expression tag	UNP P0DTC2
A	1254	HIS	-	expression tag	UNP P0DTC2
A	1255	HIS	-	expression tag	UNP P0DTC2
A	1256	HIS	-	expression tag	UNP P0DTC2
A	1257	HIS	-	expression tag	UNP P0DTC2
A	1258	SER	-	expression tag	UNP P0DTC2
A	1259	ALA	-	expression tag	UNP P0DTC2
A	1260	TRP	-	expression tag	UNP P0DTC2
A	1261	SER	-	expression tag	UNP P0DTC2
A	1262	HIS	-	expression tag	UNP P0DTC2
A	1263	PRO	-	expression tag	UNP P0DTC2
A	1264	GLN	-	expression tag	UNP P0DTC2
A	1265	PHE	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1266	GLU	-	expression tag	UNP P0DTC2
A	1267	LYS	-	expression tag	UNP P0DTC2
A	1268	GLY	-	expression tag	UNP P0DTC2
A	1269	GLY	-	expression tag	UNP P0DTC2
A	1270	GLY	-	expression tag	UNP P0DTC2
A	1271	SER	-	expression tag	UNP P0DTC2
A	1272	GLY	-	expression tag	UNP P0DTC2
A	1273	GLY	-	expression tag	UNP P0DTC2
A	1274	GLY	-	expression tag	UNP P0DTC2
A	1275	GLY	-	expression tag	UNP P0DTC2
A	1276	SER	-	expression tag	UNP P0DTC2
A	1277	GLY	-	expression tag	UNP P0DTC2
A	1278	GLY	-	expression tag	UNP P0DTC2
A	1279	SER	-	expression tag	UNP P0DTC2
A	1280	ALA	-	expression tag	UNP P0DTC2
A	1281	TRP	-	expression tag	UNP P0DTC2
A	1282	SER	-	expression tag	UNP P0DTC2
A	1283	HIS	-	expression tag	UNP P0DTC2
A	1284	PRO	-	expression tag	UNP P0DTC2
A	1285	GLN	-	expression tag	UNP P0DTC2
A	1286	PHE	-	expression tag	UNP P0DTC2
A	1287	GLU	-	expression tag	UNP P0DTC2
A	1288	LYS	-	expression tag	UNP P0DTC2
E	682	GLY	ARG	conflict	UNP P0DTC2
E	683	SER	ARG	conflict	UNP P0DTC2
E	685	SER	ARG	conflict	UNP P0DTC2
E	817	PRO	PHE	conflict	UNP P0DTC2
E	892	PRO	ALA	conflict	UNP P0DTC2
E	899	PRO	ALA	conflict	UNP P0DTC2
E	942	PRO	ALA	conflict	UNP P0DTC2
E	986	PRO	LYS	conflict	UNP P0DTC2
E	987	PRO	VAL	conflict	UNP P0DTC2
E	1209	GLY	-	expression tag	UNP P0DTC2
E	1210	SER	-	expression tag	UNP P0DTC2
E	1211	GLY	-	expression tag	UNP P0DTC2
E	1212	TYR	-	expression tag	UNP P0DTC2
E	1213	ILE	-	expression tag	UNP P0DTC2
E	1214	PRO	-	expression tag	UNP P0DTC2
E	1215	GLU	-	expression tag	UNP P0DTC2
E	1216	ALA	-	expression tag	UNP P0DTC2
E	1217	PRO	-	expression tag	UNP P0DTC2
E	1218	ARG	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
E	1219	ASP	-	expression tag	UNP P0DTC2
E	1220	GLY	-	expression tag	UNP P0DTC2
E	1221	GLN	-	expression tag	UNP P0DTC2
E	1222	ALA	-	expression tag	UNP P0DTC2
E	1223	TYR	-	expression tag	UNP P0DTC2
E	1224	VAL	-	expression tag	UNP P0DTC2
E	1225	ARG	-	expression tag	UNP P0DTC2
E	1226	LYS	-	expression tag	UNP P0DTC2
E	1227	ASP	-	expression tag	UNP P0DTC2
E	1228	GLY	-	expression tag	UNP P0DTC2
E	1229	GLU	-	expression tag	UNP P0DTC2
E	1230	TRP	-	expression tag	UNP P0DTC2
E	1231	VAL	-	expression tag	UNP P0DTC2
E	1232	LEU	-	expression tag	UNP P0DTC2
E	1233	LEU	-	expression tag	UNP P0DTC2
E	1234	SER	-	expression tag	UNP P0DTC2
E	1235	THR	-	expression tag	UNP P0DTC2
E	1236	PHE	-	expression tag	UNP P0DTC2
E	1237	LEU	-	expression tag	UNP P0DTC2
E	1238	GLY	-	expression tag	UNP P0DTC2
E	1239	ARG	-	expression tag	UNP P0DTC2
E	1240	SER	-	expression tag	UNP P0DTC2
E	1241	LEU	-	expression tag	UNP P0DTC2
E	1242	GLU	-	expression tag	UNP P0DTC2
E	1243	VAL	-	expression tag	UNP P0DTC2
E	1244	LEU	-	expression tag	UNP P0DTC2
E	1245	PHE	-	expression tag	UNP P0DTC2
E	1246	GLN	-	expression tag	UNP P0DTC2
E	1247	GLY	-	expression tag	UNP P0DTC2
E	1248	PRO	-	expression tag	UNP P0DTC2
E	1249	GLY	-	expression tag	UNP P0DTC2
E	1250	HIS	-	expression tag	UNP P0DTC2
E	1251	HIS	-	expression tag	UNP P0DTC2
E	1252	HIS	-	expression tag	UNP P0DTC2
E	1253	HIS	-	expression tag	UNP P0DTC2
E	1254	HIS	-	expression tag	UNP P0DTC2
E	1255	HIS	-	expression tag	UNP P0DTC2
E	1256	HIS	-	expression tag	UNP P0DTC2
E	1257	HIS	-	expression tag	UNP P0DTC2
E	1258	SER	-	expression tag	UNP P0DTC2
E	1259	ALA	-	expression tag	UNP P0DTC2
E	1260	TRP	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
E	1261	SER	-	expression tag	UNP P0DTC2
E	1262	HIS	-	expression tag	UNP P0DTC2
E	1263	PRO	-	expression tag	UNP P0DTC2
E	1264	GLN	-	expression tag	UNP P0DTC2
E	1265	PHE	-	expression tag	UNP P0DTC2
E	1266	GLU	-	expression tag	UNP P0DTC2
E	1267	LYS	-	expression tag	UNP P0DTC2
E	1268	GLY	-	expression tag	UNP P0DTC2
E	1269	GLY	-	expression tag	UNP P0DTC2
E	1270	GLY	-	expression tag	UNP P0DTC2
E	1271	SER	-	expression tag	UNP P0DTC2
E	1272	GLY	-	expression tag	UNP P0DTC2
E	1273	GLY	-	expression tag	UNP P0DTC2
E	1274	GLY	-	expression tag	UNP P0DTC2
E	1275	GLY	-	expression tag	UNP P0DTC2
E	1276	SER	-	expression tag	UNP P0DTC2
E	1277	GLY	-	expression tag	UNP P0DTC2
E	1278	GLY	-	expression tag	UNP P0DTC2
E	1279	SER	-	expression tag	UNP P0DTC2
E	1280	ALA	-	expression tag	UNP P0DTC2
E	1281	TRP	-	expression tag	UNP P0DTC2
E	1282	SER	-	expression tag	UNP P0DTC2
E	1283	HIS	-	expression tag	UNP P0DTC2
E	1284	PRO	-	expression tag	UNP P0DTC2
E	1285	GLN	-	expression tag	UNP P0DTC2
E	1286	PHE	-	expression tag	UNP P0DTC2
E	1287	GLU	-	expression tag	UNP P0DTC2
E	1288	LYS	-	expression tag	UNP P0DTC2
G	682	GLY	ARG	conflict	UNP P0DTC2
G	683	SER	ARG	conflict	UNP P0DTC2
G	685	SER	ARG	conflict	UNP P0DTC2
G	817	PRO	PHE	conflict	UNP P0DTC2
G	892	PRO	ALA	conflict	UNP P0DTC2
G	899	PRO	ALA	conflict	UNP P0DTC2
G	942	PRO	ALA	conflict	UNP P0DTC2
G	986	PRO	LYS	conflict	UNP P0DTC2
G	987	PRO	VAL	conflict	UNP P0DTC2
G	1209	GLY	-	expression tag	UNP P0DTC2
G	1210	SER	-	expression tag	UNP P0DTC2
G	1211	GLY	-	expression tag	UNP P0DTC2
G	1212	TYR	-	expression tag	UNP P0DTC2
G	1213	ILE	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
G	1214	PRO	-	expression tag	UNP P0DTC2
G	1215	GLU	-	expression tag	UNP P0DTC2
G	1216	ALA	-	expression tag	UNP P0DTC2
G	1217	PRO	-	expression tag	UNP P0DTC2
G	1218	ARG	-	expression tag	UNP P0DTC2
G	1219	ASP	-	expression tag	UNP P0DTC2
G	1220	GLY	-	expression tag	UNP P0DTC2
G	1221	GLN	-	expression tag	UNP P0DTC2
G	1222	ALA	-	expression tag	UNP P0DTC2
G	1223	TYR	-	expression tag	UNP P0DTC2
G	1224	VAL	-	expression tag	UNP P0DTC2
G	1225	ARG	-	expression tag	UNP P0DTC2
G	1226	LYS	-	expression tag	UNP P0DTC2
G	1227	ASP	-	expression tag	UNP P0DTC2
G	1228	GLY	-	expression tag	UNP P0DTC2
G	1229	GLU	-	expression tag	UNP P0DTC2
G	1230	TRP	-	expression tag	UNP P0DTC2
G	1231	VAL	-	expression tag	UNP P0DTC2
G	1232	LEU	-	expression tag	UNP P0DTC2
G	1233	LEU	-	expression tag	UNP P0DTC2
G	1234	SER	-	expression tag	UNP P0DTC2
G	1235	THR	-	expression tag	UNP P0DTC2
G	1236	PHE	-	expression tag	UNP P0DTC2
G	1237	LEU	-	expression tag	UNP P0DTC2
G	1238	GLY	-	expression tag	UNP P0DTC2
G	1239	ARG	-	expression tag	UNP P0DTC2
G	1240	SER	-	expression tag	UNP P0DTC2
G	1241	LEU	-	expression tag	UNP P0DTC2
G	1242	GLU	-	expression tag	UNP P0DTC2
G	1243	VAL	-	expression tag	UNP P0DTC2
G	1244	LEU	-	expression tag	UNP P0DTC2
G	1245	PHE	-	expression tag	UNP P0DTC2
G	1246	GLN	-	expression tag	UNP P0DTC2
G	1247	GLY	-	expression tag	UNP P0DTC2
G	1248	PRO	-	expression tag	UNP P0DTC2
G	1249	GLY	-	expression tag	UNP P0DTC2
G	1250	HIS	-	expression tag	UNP P0DTC2
G	1251	HIS	-	expression tag	UNP P0DTC2
G	1252	HIS	-	expression tag	UNP P0DTC2
G	1253	HIS	-	expression tag	UNP P0DTC2
G	1254	HIS	-	expression tag	UNP P0DTC2
G	1255	HIS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
G	1256	HIS	-	expression tag	UNP P0DTC2
G	1257	HIS	-	expression tag	UNP P0DTC2
G	1258	SER	-	expression tag	UNP P0DTC2
G	1259	ALA	-	expression tag	UNP P0DTC2
G	1260	TRP	-	expression tag	UNP P0DTC2
G	1261	SER	-	expression tag	UNP P0DTC2
G	1262	HIS	-	expression tag	UNP P0DTC2
G	1263	PRO	-	expression tag	UNP P0DTC2
G	1264	GLN	-	expression tag	UNP P0DTC2
G	1265	PHE	-	expression tag	UNP P0DTC2
G	1266	GLU	-	expression tag	UNP P0DTC2
G	1267	LYS	-	expression tag	UNP P0DTC2
G	1268	GLY	-	expression tag	UNP P0DTC2
G	1269	GLY	-	expression tag	UNP P0DTC2
G	1270	GLY	-	expression tag	UNP P0DTC2
G	1271	SER	-	expression tag	UNP P0DTC2
G	1272	GLY	-	expression tag	UNP P0DTC2
G	1273	GLY	-	expression tag	UNP P0DTC2
G	1274	GLY	-	expression tag	UNP P0DTC2
G	1275	GLY	-	expression tag	UNP P0DTC2
G	1276	SER	-	expression tag	UNP P0DTC2
G	1277	GLY	-	expression tag	UNP P0DTC2
G	1278	GLY	-	expression tag	UNP P0DTC2
G	1279	SER	-	expression tag	UNP P0DTC2
G	1280	ALA	-	expression tag	UNP P0DTC2
G	1281	TRP	-	expression tag	UNP P0DTC2
G	1282	SER	-	expression tag	UNP P0DTC2
G	1283	HIS	-	expression tag	UNP P0DTC2
G	1284	PRO	-	expression tag	UNP P0DTC2
G	1285	GLN	-	expression tag	UNP P0DTC2
G	1286	PHE	-	expression tag	UNP P0DTC2
G	1287	GLU	-	expression tag	UNP P0DTC2
G	1288	LYS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called Fab P5C3 Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	108	Total	C	N	O	S	0	0
			826	519	142	163	2		
2	D	108	Total	C	N	O	S	0	0
			826	519	142	163	2		
2	H	108	Total	C	N	O	S	0	0
			826	519	142	163	2		

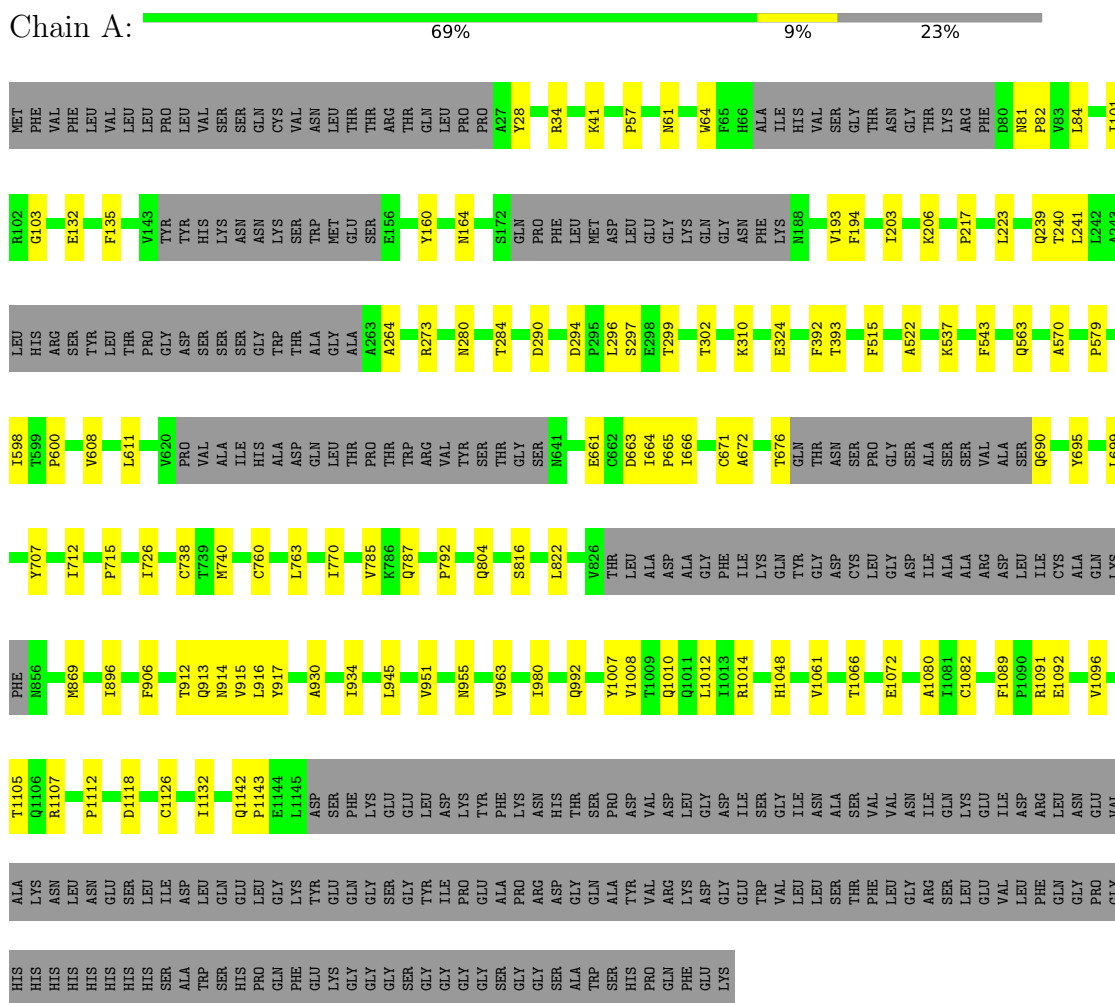
- Molecule 3 is a protein called Fab P5C3 HEAVY Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	123	Total	C	N	O	S	0	0
			932	579	158	187	8		
3	F	123	Total	C	N	O	S	0	0
			932	579	158	187	8		
3	I	123	Total	C	N	O	S	0	0
			932	579	158	187	8		

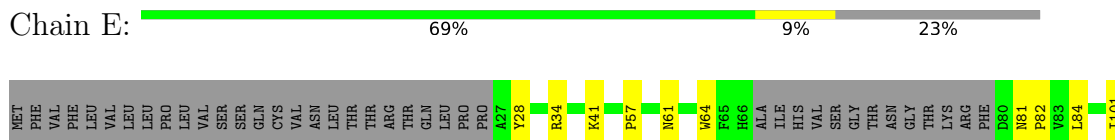
3 Residue-property plots

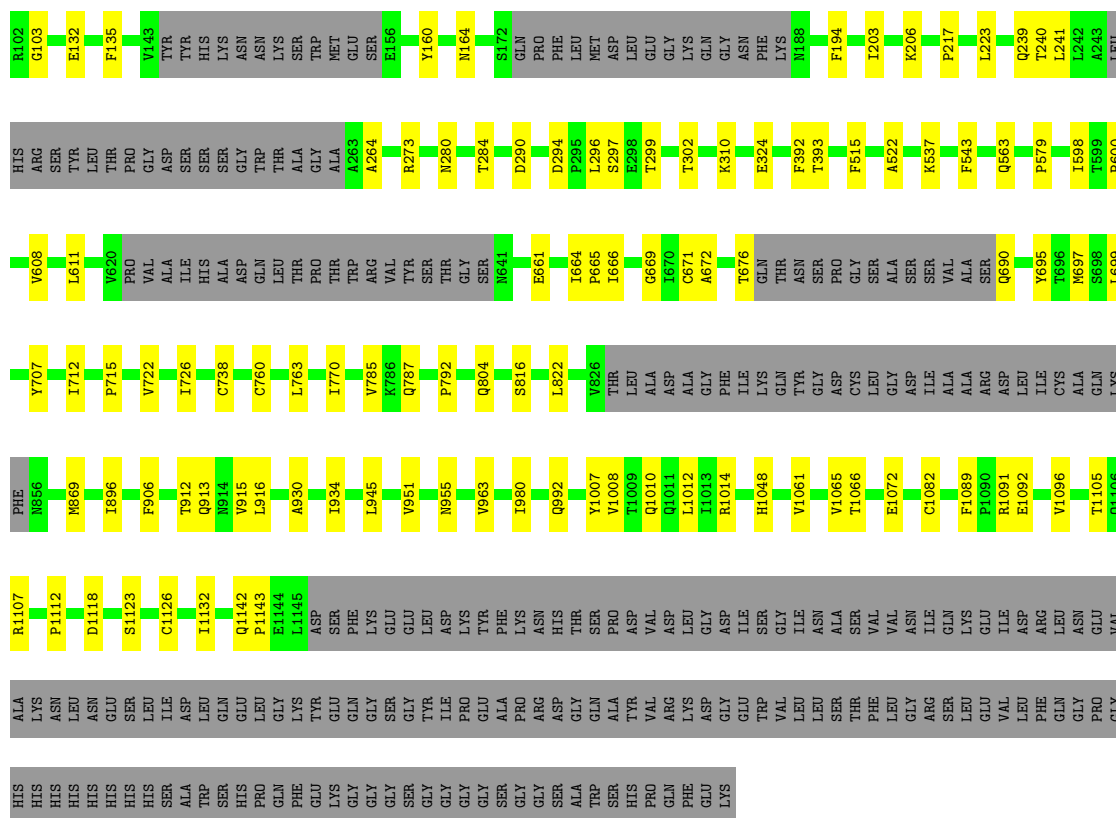
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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

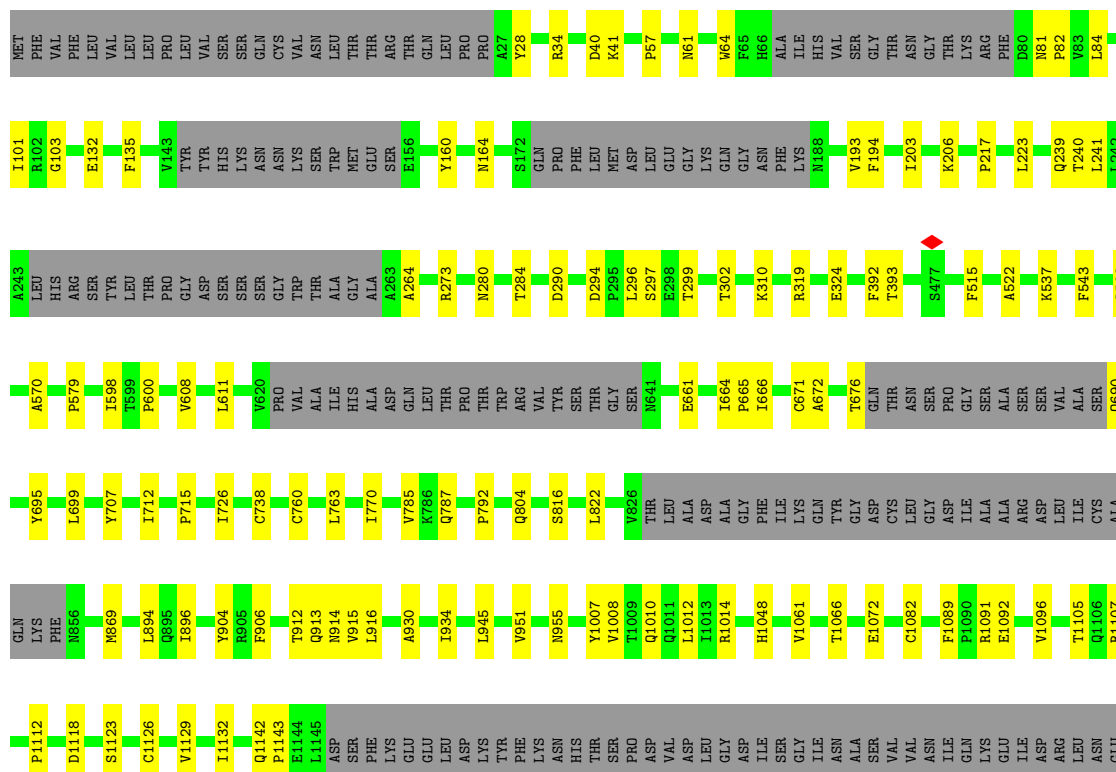


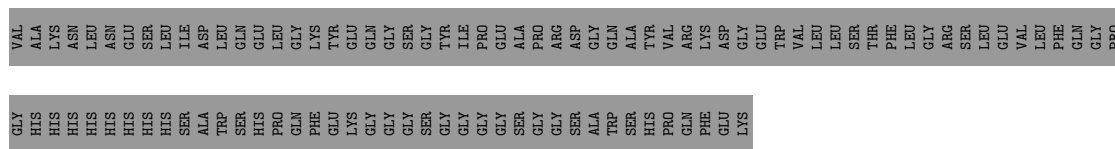
• Molecule 1: Spike glycoprotein



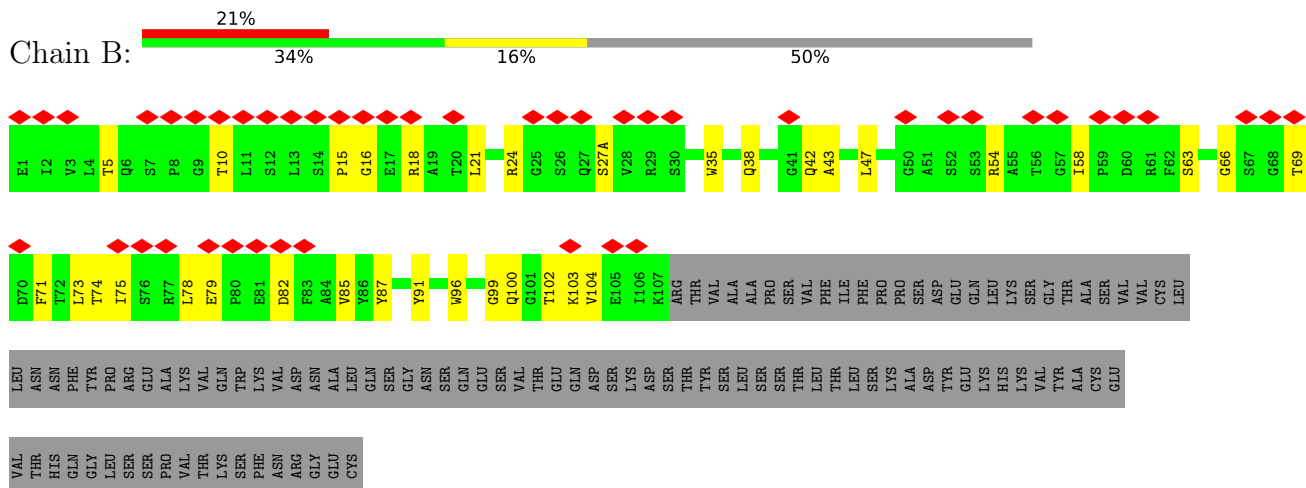


● Molecule 1: Spike glycoprotein

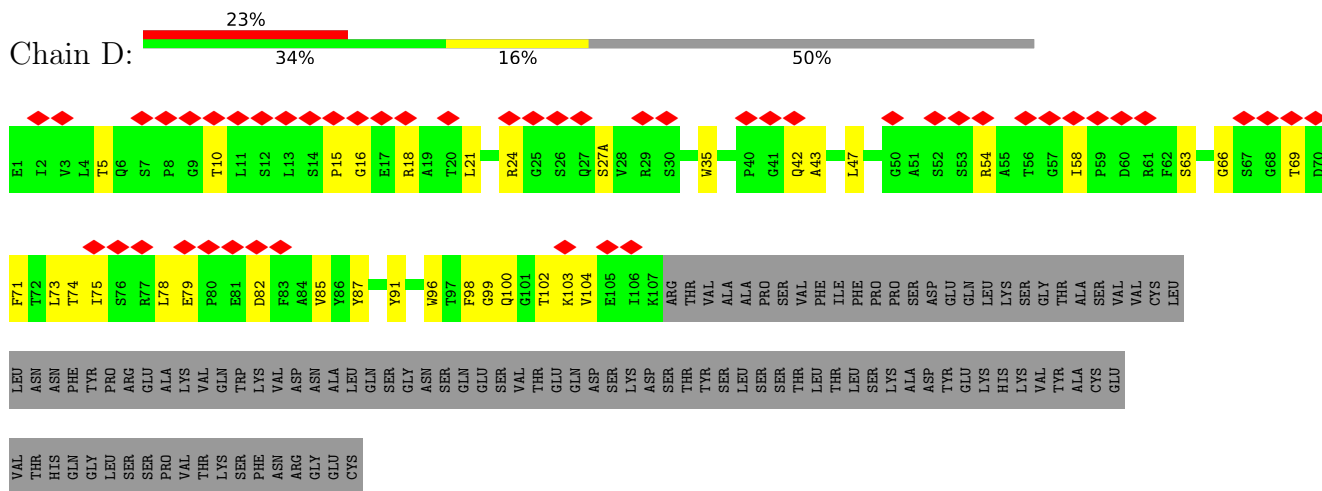




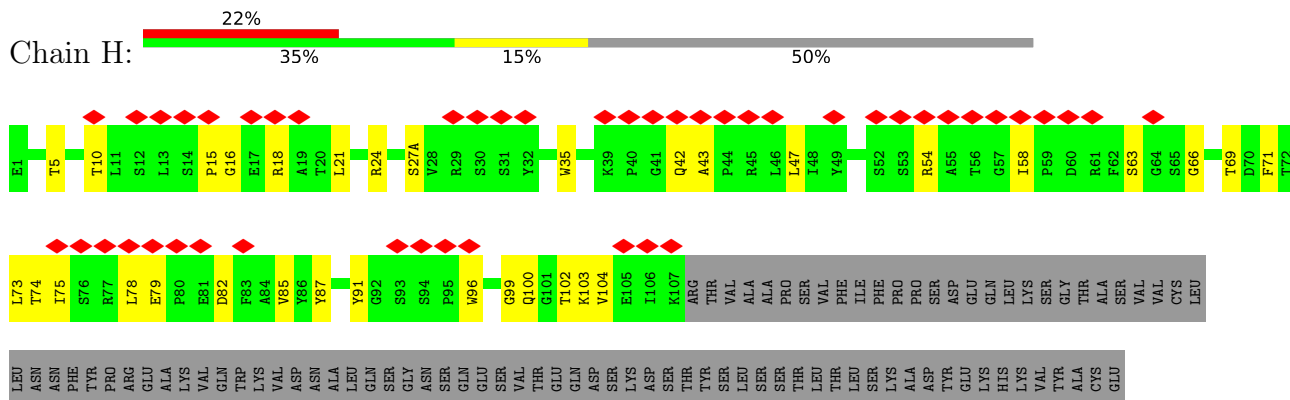
- Molecule 2: Fab P5C3 Light Chain



- Molecule 2: Fab P5C3 Light Chain

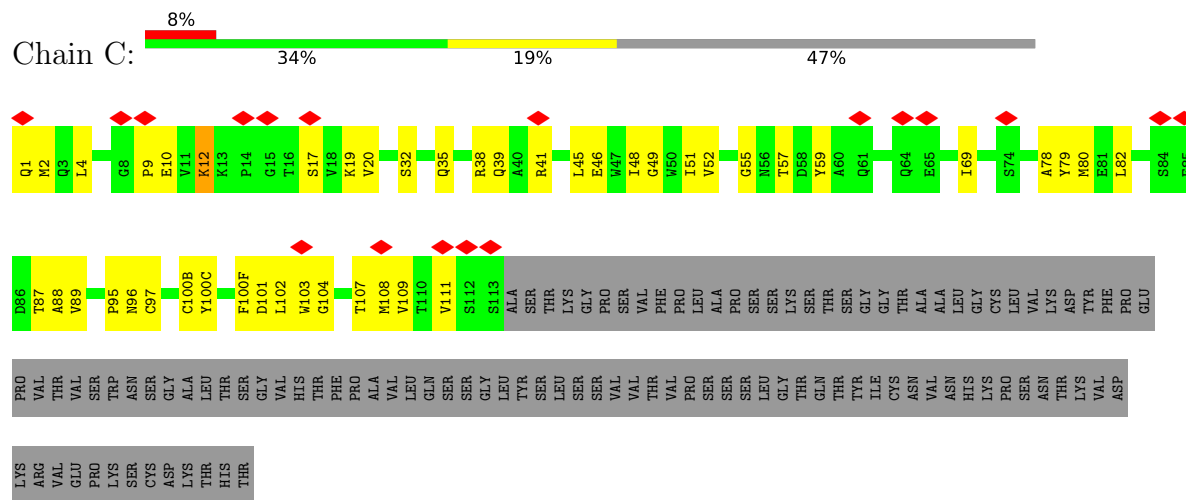


- Molecule 2: Fab P5C3 Light Chain

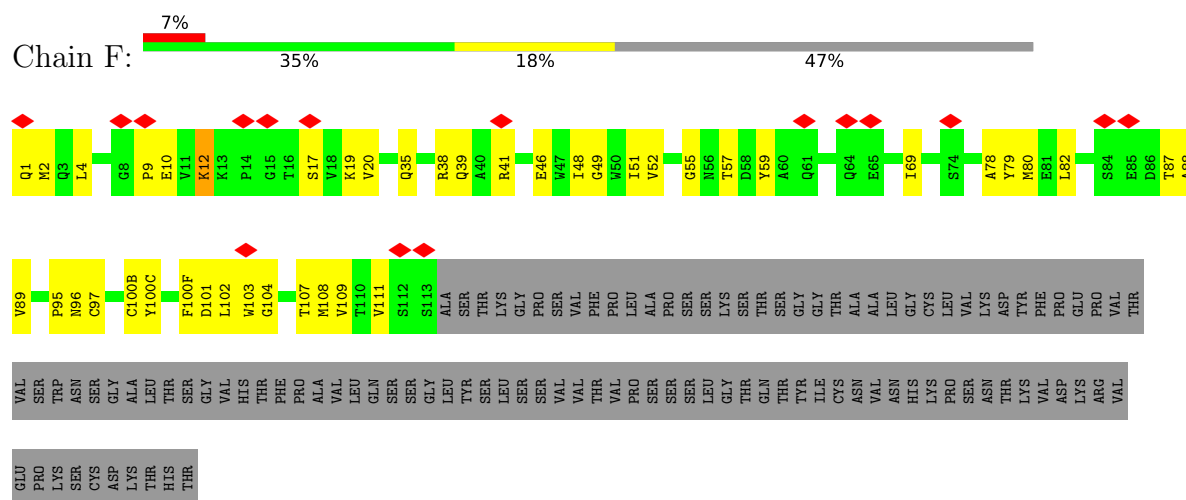


VAL
THR
HIS
GLN
GLY
LEU
SER
SER
PRO
VAL
THR
LYS
SER
PHE
ASN
ARG
GLY
CYS

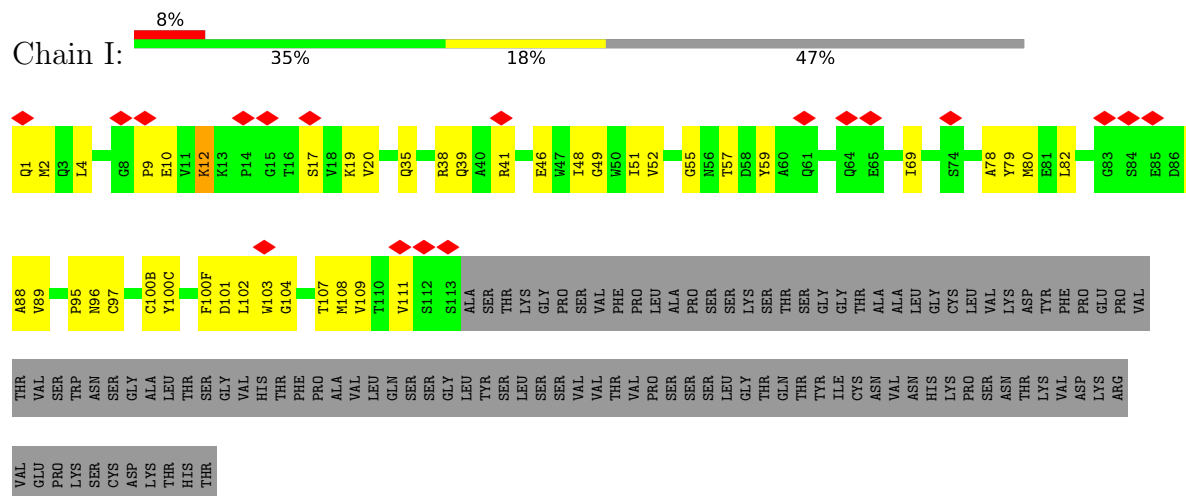
• Molecule 3: Fab P5C3 HEAVY Chain



• Molecule 3: Fab P5C3 HEAVY Chain



• Molecule 3: Fab P5C3 HEAVY Chain



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	120000	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 ($e^-/\text{\AA}^2$)	38	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	105000	Depositor
Image detector	GATAN K2 IS (4k x 4k)	Depositor
Maximum map value	3.457	Depositor
Minimum map value	-1.320	Depositor
Average map value	-0.004	Depositor
Map value standard deviation	0.059	Depositor
Recommended contour level	0.3	Depositor
Map size (Å)	492.0, 492.0, 492.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.64, 1.64, 1.64	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	A	0.26	0/7222	0.46	0/9891
1	E	0.26	0/7222	0.46	0/9891
1	G	0.26	0/7222	0.46	0/9891
2	B	0.29	0/846	0.56	0/1147
2	D	0.29	0/846	0.56	0/1147
2	H	0.29	0/846	0.56	0/1147
3	C	0.27	0/951	0.59	0/1289
3	F	0.27	0/951	0.59	0/1289
3	I	0.27	0/951	0.59	0/1289
All	All	0.26	0/27057	0.48	0/36981

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7064	0	6245	73	0
1	E	7064	0	6245	71	0
1	G	7064	0	6245	73	0
2	B	826	0	802	21	0
2	D	826	0	802	24	0
2	H	826	0	802	22	0
3	C	932	0	887	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	932	0	887	38	0
3	I	932	0	887	39	0
All	All	26466	0	23802	373	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 (373) 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)
3:I:10:GLU:HB3	3:I:108:MET:O	1.66	0.95
3:C:10:GLU:HB3	3:C:108:MET:O	1.67	0.95
3:F:10:GLU:HB3	3:F:108:MET:O	1.67	0.95
3:I:17:SER:HA	3:I:82:LEU:O	1.81	0.81
3:C:17:SER:HA	3:C:82:LEU:O	1.81	0.80
3:F:17:SER:HA	3:F:82:LEU:O	1.81	0.80
1:E:132:GLU:HB2	1:E:164:ASN:O	1.82	0.79
1:G:132:GLU:HB2	1:G:164:ASN:O	1.82	0.79
1:A:132:GLU:HB2	1:A:164:ASN:O	1.82	0.78
3:C:45:LEU:HD21	2:D:87:TYR:OH	1.85	0.77
3:F:2:MET:HG2	3:F:102:LEU:HD12	1.68	0.76
3:C:2:MET:HG2	3:C:102:LEU:HD12	1.68	0.74
3:I:2:MET:HG2	3:I:102:LEU:HD12	1.68	0.74
3:I:9:PRO:HA	3:I:108:MET:HB2	1.70	0.73
2:B:38:GLN:NE2	3:F:39:GLN:OE1	2.20	0.73
3:F:9:PRO:HA	3:F:108:MET:HB2	1.70	0.73
3:C:9:PRO:HA	3:C:108:MET:HB2	1.70	0.72
2:B:47:LEU:HD12	2:B:58:ILE:HD12	1.71	0.71
2:D:47:LEU:HD12	2:D:58:ILE:HD12	1.72	0.71
2:H:47:LEU:HD12	2:H:58:ILE:HD12	1.72	0.71
1:E:103:GLY:HA3	1:E:241:LEU:HB2	1.74	0.70
1:A:103:GLY:HA3	1:A:241:LEU:HB2	1.74	0.69
1:G:103:GLY:HA3	1:G:241:LEU:HB2	1.74	0.69
2:H:5:THR:HG21	3:I:100(F):PHE:HD2	1.59	0.68
2:H:5:THR:HG21	3:I:100(F):PHE:CD2	2.30	0.66
1:E:707:TYR:HB3	1:G:792:PRO:HG3	1.78	0.66
3:C:87:THR:HB	3:C:111:VAL:HG12	1.78	0.65
3:C:4:LEU:HB2	3:C:104:GLY:HA2	1.79	0.65
3:I:87:THR:HB	3:I:111:VAL:HG12	1.78	0.65
1:A:763:LEU:HD12	1:A:1008:VAL:HG21	1.79	0.64
3:C:69:ILE:HD11	3:C:78:ALA:HB1	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:4:LEU:HB2	3:F:104:GLY:HA2	1.79	0.64
3:F:69:ILE:HD11	3:F:78:ALA:HB1	1.79	0.64
3:F:87:THR:HB	3:F:111:VAL:HG12	1.78	0.64
3:I:38:ARG:NE	3:I:46:GLU:OE2	2.30	0.64
3:I:69:ILE:HD11	3:I:78:ALA:HB1	1.79	0.64
3:C:38:ARG:NE	3:C:46:GLU:OE2	2.30	0.64
1:G:763:LEU:HD12	1:G:1008:VAL:HG21	1.79	0.64
1:A:1089:PHE:HB3	1:E:913:GLN:HE21	1.63	0.64
1:G:726:ILE:HG12	1:G:1061:VAL:HG22	1.80	0.63
1:A:726:ILE:HG12	1:A:1061:VAL:HG22	1.80	0.63
3:F:38:ARG:NE	3:F:46:GLU:OE2	2.30	0.63
3:I:4:LEU:HB2	3:I:104:GLY:HA2	1.79	0.63
1:E:763:LEU:HD12	1:E:1008:VAL:HG21	1.79	0.62
3:I:49:GLY:HA3	3:I:59:TYR:HD1	1.64	0.62
1:E:726:ILE:HG12	1:E:1061:VAL:HG22	1.80	0.62
3:F:49:GLY:HA3	3:F:59:TYR:HD1	1.64	0.62
1:E:699:LEU:HD11	1:G:869:MET:HB2	1.81	0.62
1:E:81:ASN:O	1:E:239:GLN:NE2	2.33	0.62
1:E:324:GLU:OE2	1:E:537:LYS:NZ	2.33	0.61
3:C:49:GLY:HA3	3:C:59:TYR:HD1	1.64	0.61
1:A:869:MET:HB2	1:G:699:LEU:HD11	1.82	0.61
3:C:96:ASN:ND2	3:C:101:ASP:HB2	2.16	0.61
3:F:38:ARG:HB3	3:F:48:ILE:HD11	1.83	0.61
1:G:324:GLU:OE2	1:G:537:LYS:NZ	2.33	0.60
1:A:324:GLU:OE2	1:A:537:LYS:NZ	2.33	0.60
3:I:38:ARG:HB3	3:I:48:ILE:HD11	1.83	0.60
1:A:64:TRP:HE1	1:A:264:ALA:HB1	1.67	0.60
2:H:35:TRP:HA	2:H:87:TYR:O	2.02	0.60
3:I:96:ASN:ND2	3:I:102:LEU:HG	2.17	0.60
1:E:1089:PHE:HB3	1:G:913:GLN:HE21	1.67	0.60
3:F:96:ASN:ND2	3:F:101:ASP:HB2	2.16	0.60
1:E:1092:GLU:O	1:E:1107:ARG:NH1	2.35	0.60
2:B:35:TRP:HA	2:B:87:TYR:O	2.02	0.60
3:F:39:GLN:O	3:F:88:ALA:HA	2.02	0.60
1:A:1092:GLU:O	1:A:1107:ARG:NH1	2.35	0.60
3:C:96:ASN:ND2	3:C:102:LEU:HG	2.17	0.59
2:D:35:TRP:HA	2:D:87:TYR:O	2.02	0.59
1:A:913:GLN:HE21	1:G:1089:PHE:HB3	1.66	0.59
3:C:38:ARG:HB3	3:C:48:ILE:HD11	1.83	0.59
3:F:96:ASN:ND2	3:F:102:LEU:HG	2.17	0.59
1:G:64:TRP:HE1	1:G:264:ALA:HB1	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:96:ASN:ND2	3:I:101:ASP:HB2	2.16	0.59
2:B:63:SER:OG	2:B:74:THR:OG1	2.20	0.59
1:E:64:TRP:HE1	1:E:264:ALA:HB1	1.67	0.59
3:C:39:GLN:O	3:C:88:ALA:HA	2.02	0.59
2:H:63:SER:OG	2:H:74:THR:OG1	2.20	0.59
2:D:63:SER:OG	2:D:74:THR:OG1	2.20	0.59
1:G:951:VAL:O	1:G:955:ASN:ND2	2.36	0.59
1:G:1092:GLU:O	1:G:1107:ARG:NH1	2.35	0.59
1:A:951:VAL:O	1:A:955:ASN:ND2	2.36	0.58
3:I:39:GLN:O	3:I:88:ALA:HA	2.02	0.58
1:A:41:LYS:HB3	1:G:563:GLN:HA	1.85	0.58
3:F:12:LYS:HE3	3:F:111:VAL:HB	1.86	0.58
1:G:280:ASN:ND2	1:G:284:THR:OG1	2.37	0.58
3:F:1:GLN:HG3	3:F:102:LEU:HD11	1.86	0.57
2:D:16:GLY:H	2:D:78:LEU:HB3	1.69	0.57
2:D:54:ARG:HH12	2:D:63:SER:HA	1.69	0.57
2:B:16:GLY:H	2:B:78:LEU:HB3	1.69	0.57
3:C:1:GLN:HG3	3:C:102:LEU:HD11	1.86	0.57
2:H:18:ARG:HH21	2:H:74:THR:HG21	1.70	0.57
1:E:951:VAL:O	1:E:955:ASN:ND2	2.36	0.57
3:I:12:LYS:HE3	3:I:111:VAL:HB	1.86	0.57
1:A:81:ASN:O	1:A:239:GLN:NE2	2.32	0.57
2:H:54:ARG:HH12	2:H:63:SER:HA	1.69	0.57
1:A:280:ASN:ND2	1:A:284:THR:OG1	2.37	0.57
2:D:18:ARG:HH21	2:D:74:THR:HG21	1.70	0.57
1:E:280:ASN:ND2	1:E:284:THR:OG1	2.37	0.57
2:B:54:ARG:HH12	2:B:63:SER:HA	1.69	0.57
3:C:12:LYS:HE3	3:C:111:VAL:HB	1.86	0.57
1:G:770:ILE:HD11	1:G:1012:LEU:HD23	1.87	0.57
1:A:699:LEU:HD11	1:E:869:MET:HB2	1.86	0.56
2:H:16:GLY:H	2:H:78:LEU:HB3	1.69	0.56
2:B:18:ARG:HH21	2:B:74:THR:HG21	1.70	0.56
1:G:81:ASN:O	1:G:239:GLN:NE2	2.32	0.56
3:I:1:GLN:HG3	3:I:102:LEU:HD11	1.86	0.56
1:A:770:ILE:HD11	1:A:1012:LEU:HD23	1.87	0.56
3:C:19:LYS:HZ1	3:C:79:TYR:HD1	1.53	0.56
1:G:393:THR:HA	1:G:522:ALA:HA	1.88	0.56
3:C:103:TRP:CZ3	2:D:42:GLN:HG3	2.41	0.56
1:A:792:PRO:HG3	1:G:707:TYR:HB3	1.88	0.56
1:A:804:GLN:O	1:A:816:SER:OG	2.24	0.56
3:I:52:VAL:O	3:I:55:GLY:N	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:804:GLN:O	1:G:816:SER:OG	2.24	0.56
3:C:52:VAL:O	3:C:55:GLY:N	2.39	0.55
3:F:52:VAL:O	3:F:55:GLY:N	2.39	0.55
1:E:770:ILE:HD11	1:E:1012:LEU:HD23	1.87	0.55
1:A:393:THR:HA	1:A:522:ALA:HA	1.88	0.55
2:D:15:PRO:HA	2:D:78:LEU:HB3	1.89	0.55
1:E:804:GLN:O	1:E:816:SER:OG	2.24	0.55
2:H:15:PRO:HA	2:H:78:LEU:HB3	1.89	0.54
2:B:15:PRO:HA	2:B:78:LEU:HB3	1.89	0.54
1:E:393:THR:HA	1:E:522:ALA:HA	1.88	0.54
3:I:20:VAL:HG13	3:I:107:THR:HG21	1.90	0.54
1:A:707:TYR:HB3	1:E:792:PRO:HG3	1.90	0.54
1:A:896:ILE:HD12	1:G:712:ILE:HG22	1.89	0.54
1:A:712:ILE:HD11	1:A:1096:VAL:HG12	1.90	0.54
3:I:41:ARG:HD2	3:I:41:ARG:O	2.08	0.54
1:E:712:ILE:HD11	1:E:1096:VAL:HG12	1.90	0.53
3:F:41:ARG:HD2	3:F:41:ARG:O	2.08	0.53
3:C:20:VAL:HG13	3:C:107:THR:HG21	1.90	0.53
3:F:19:LYS:HZ1	3:F:79:TYR:HD1	1.57	0.53
1:A:666:ILE:HD11	1:A:672:ALA:HB2	1.91	0.53
3:C:41:ARG:HD2	3:C:41:ARG:O	2.08	0.53
3:I:95:PRO:HA	3:I:100(F):PHE:CD1	2.44	0.53
1:G:712:ILE:HD11	1:G:1096:VAL:HG12	1.90	0.53
3:C:95:PRO:HA	3:C:100(F):PHE:CD1	2.44	0.53
1:G:392:PHE:CG	1:G:515:PHE:HB3	2.44	0.52
1:E:392:PHE:CG	1:E:515:PHE:HB3	2.44	0.52
3:F:20:VAL:HG13	3:F:107:THR:HG21	1.90	0.52
3:F:95:PRO:HA	3:F:100(F):PHE:CD1	2.44	0.52
3:F:35:GLN:NE2	3:F:100(F):PHE:HE2	2.08	0.52
1:G:1010:GLN:OE1	1:G:1014:ARG:NH1	2.39	0.52
2:H:66:GLY:HA3	2:H:71:PHE:HA	1.91	0.52
1:A:1091:ARG:NH1	1:A:1118:ASP:O	2.43	0.52
3:C:35:GLN:NE2	3:C:100(F):PHE:HE2	2.08	0.52
3:F:96:ASN:HD22	3:F:101:ASP:HB2	1.75	0.52
1:A:392:PHE:CG	1:A:515:PHE:HB3	2.44	0.51
1:A:1010:GLN:OE1	1:A:1014:ARG:NH1	2.39	0.51
2:D:66:GLY:HA3	2:D:71:PHE:HA	1.91	0.51
1:G:906:PHE:CD2	1:G:916:LEU:HB2	2.45	0.51
3:C:103:TRP:HZ3	2:D:42:GLN:HG3	1.74	0.51
1:E:1091:ARG:NH1	1:E:1118:ASP:O	2.43	0.51
3:I:35:GLN:NE2	3:I:100(F):PHE:HE2	2.08	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:906:PHE:CD2	1:E:916:LEU:HB2	2.46	0.51
1:G:666:ILE:HD11	1:G:672:ALA:HB2	1.91	0.51
1:A:101:ILE:HD11	1:A:240:THR:HB	1.92	0.51
1:A:294:ASP:OD1	1:A:297:SER:N	2.39	0.51
1:E:101:ILE:HD11	1:E:240:THR:HB	1.92	0.51
1:G:101:ILE:HD11	1:G:240:THR:HB	1.92	0.51
3:C:89:VAL:HA	3:C:107:THR:O	2.11	0.51
1:G:1091:ARG:NH1	1:G:1118:ASP:O	2.43	0.51
3:I:38:ARG:HA	3:I:89:VAL:O	2.11	0.51
1:A:661:GLU:O	1:A:695:TYR:OH	2.27	0.51
3:F:89:VAL:HA	3:F:107:THR:O	2.11	0.51
1:A:906:PHE:CD2	1:A:916:LEU:HB2	2.46	0.50
1:E:666:ILE:HD11	1:E:672:ALA:HB2	1.91	0.50
2:B:66:GLY:HA3	2:B:71:PHE:HA	1.91	0.50
3:C:96:ASN:HD22	3:C:101:ASP:HB2	1.75	0.50
3:C:88:ALA:O	3:C:109:VAL:HB	2.11	0.50
1:E:661:GLU:O	1:E:695:TYR:OH	2.27	0.50
3:I:17:SER:CA	3:I:82:LEU:O	2.58	0.50
1:A:563:GLN:HA	1:E:41:LYS:HB3	1.93	0.50
1:G:715:PRO:HA	1:G:1072:GLU:HA	1.94	0.50
3:C:17:SER:CA	3:C:82:LEU:O	2.58	0.50
3:F:38:ARG:HA	3:F:89:VAL:O	2.11	0.50
1:A:1126:CYS:HB3	1:A:1132:ILE:HD13	1.94	0.50
1:E:1010:GLN:OE1	1:E:1014:ARG:NH1	2.39	0.49
3:F:88:ALA:O	3:F:109:VAL:HB	2.12	0.49
1:E:715:PRO:HA	1:E:1072:GLU:HA	1.94	0.49
1:A:715:PRO:HA	1:A:1072:GLU:HA	1.93	0.49
2:B:10:THR:HA	2:B:103:LYS:HD2	1.95	0.49
1:E:294:ASP:OD1	1:E:297:SER:N	2.39	0.49
3:I:89:VAL:HA	3:I:107:THR:O	2.11	0.49
1:E:563:GLN:HA	1:G:41:LYS:HB3	1.94	0.49
2:H:10:THR:HA	2:H:103:LYS:HD2	1.95	0.49
2:D:10:THR:HA	2:D:103:LYS:HD2	1.95	0.49
2:H:75:ILE:HG12	2:H:82:ASP:OD2	2.13	0.49
2:H:91:TYR:HB2	2:H:96:TRP:CZ2	2.48	0.49
3:C:38:ARG:HA	3:C:89:VAL:O	2.11	0.49
1:E:1126:CYS:HB3	1:E:1132:ILE:HD13	1.94	0.49
1:G:822:LEU:HD22	1:G:945:LEU:HD21	1.95	0.49
2:D:91:TYR:HB2	2:D:96:TRP:CZ2	2.48	0.48
3:I:88:ALA:O	3:I:109:VAL:HB	2.12	0.48
1:A:822:LEU:HD22	1:A:945:LEU:HD21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:51:ILE:HD13	3:C:57:THR:HG22	1.95	0.48
1:G:661:GLU:O	1:G:695:TYR:OH	2.27	0.48
1:G:1126:CYS:HB3	1:G:1132:ILE:HD13	1.94	0.48
3:F:51:ILE:HD13	3:F:57:THR:HG22	1.95	0.48
3:I:96:ASN:HD22	3:I:101:ASP:HB2	1.75	0.48
3:I:51:ILE:HD13	3:I:57:THR:HG22	1.95	0.48
1:E:135:PHE:HA	1:E:160:TYR:HA	1.96	0.48
1:G:1048:HIS:HA	1:G:1066:THR:HG22	1.95	0.48
2:B:91:TYR:HB2	2:B:96:TRP:CZ2	2.48	0.48
1:E:822:LEU:HD22	1:E:945:LEU:HD21	1.95	0.48
2:B:75:ILE:HG12	2:B:82:ASP:OD2	2.13	0.48
1:E:1105:THR:HG22	1:E:1112:PRO:HA	1.96	0.48
1:G:1105:THR:HG22	1:G:1112:PRO:HA	1.96	0.48
1:A:930:ALA:O	1:A:934:ILE:HG12	2.14	0.48
2:D:75:ILE:HG12	2:D:82:ASP:OD2	2.13	0.48
1:G:930:ALA:O	1:G:934:ILE:HG12	2.14	0.48
1:E:712:ILE:HG22	1:G:896:ILE:HD12	1.96	0.47
1:E:1048:HIS:HA	1:E:1066:THR:HG22	1.96	0.47
1:G:135:PHE:HA	1:G:160:TYR:HA	1.96	0.47
1:G:294:ASP:OD1	1:G:297:SER:N	2.39	0.47
1:A:1048:HIS:HA	1:A:1066:THR:HG22	1.95	0.47
1:E:930:ALA:O	1:E:934:ILE:HG12	2.14	0.47
1:E:1082:CYS:HB2	1:E:1126:CYS:HB3	1.64	0.47
1:G:1082:CYS:HB2	1:G:1126:CYS:HB3	1.64	0.47
3:I:19:LYS:HZ1	3:I:79:TYR:HD1	1.61	0.47
1:A:1105:THR:HG22	1:A:1112:PRO:HA	1.96	0.47
2:D:21:LEU:HD21	2:D:104:VAL:HG23	1.97	0.47
1:E:1007:TYR:O	1:E:1010:GLN:HG3	2.15	0.47
3:F:69:ILE:HD12	3:F:80:MET:HB3	1.97	0.47
2:H:21:LEU:HD21	2:H:104:VAL:HG23	1.97	0.47
1:A:135:PHE:HA	1:A:160:TYR:HA	1.96	0.47
2:B:21:LEU:HD21	2:B:104:VAL:HG23	1.97	0.47
3:F:4:LEU:HB2	3:F:103:TRP:O	2.15	0.47
1:G:1007:TYR:O	1:G:1010:GLN:HG3	2.15	0.47
1:G:1142:GLN:HB2	1:G:1143:PRO:HD3	1.97	0.47
3:C:4:LEU:HB2	3:C:103:TRP:O	2.15	0.47
1:G:310:LYS:HG3	1:G:600:PRO:HA	1.98	0.47
3:I:4:LEU:HB2	3:I:103:TRP:O	2.15	0.47
3:I:69:ILE:HD12	3:I:80:MET:HB3	1.97	0.47
1:A:57:PRO:HG3	1:A:273:ARG:HD2	1.97	0.46
3:C:10:GLU:CB	3:C:108:MET:O	2.53	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1142:GLN:HB2	1:E:1143:PRO:HD3	1.97	0.46
1:A:1082:CYS:HB2	1:A:1126:CYS:HB3	1.64	0.46
1:E:310:LYS:HG3	1:E:600:PRO:HA	1.98	0.46
1:E:34:ARG:HH21	1:E:217:PRO:HB2	1.81	0.46
1:E:82:PRO:HG2	1:E:84:LEU:HD21	1.98	0.46
1:A:310:LYS:HG3	1:A:600:PRO:HA	1.98	0.46
1:G:296:LEU:HB2	1:G:608:VAL:HG11	1.98	0.46
2:B:42:GLN:HG3	2:B:43:ALA:H	1.81	0.46
3:C:69:ILE:HD12	3:C:80:MET:HB3	1.97	0.46
2:D:42:GLN:HG3	2:D:43:ALA:H	1.81	0.46
1:E:296:LEU:HB2	1:E:608:VAL:HG11	1.98	0.46
1:G:280:ASN:OD1	1:G:284:THR:N	2.49	0.46
1:A:1007:TYR:O	1:A:1010:GLN:HG3	2.15	0.46
1:G:57:PRO:HG3	1:G:273:ARG:HD2	1.97	0.46
2:H:42:GLN:HG3	2:H:43:ALA:H	1.81	0.46
3:I:19:LYS:HZ1	3:I:79:TYR:HB3	1.81	0.46
3:I:102:LEU:HB3	3:I:103:TRP:H	1.59	0.45
2:H:27(A):SER:HA	2:H:69:THR:HG22	1.99	0.45
1:A:1142:GLN:HB2	1:A:1143:PRO:HD3	1.97	0.45
1:A:82:PRO:HG2	1:A:84:LEU:HD21	1.98	0.45
1:G:34:ARG:HH21	1:G:217:PRO:HB2	1.81	0.45
1:A:296:LEU:HB2	1:A:608:VAL:HG11	1.98	0.45
3:C:20:VAL:HB	3:C:80:MET:SD	2.57	0.45
1:E:57:PRO:HG3	1:E:273:ARG:HD2	1.97	0.45
1:E:280:ASN:OD1	1:E:284:THR:N	2.49	0.45
3:I:20:VAL:HB	3:I:80:MET:SD	2.57	0.45
2:D:87:TYR:HD1	2:D:99:GLY:HA3	1.82	0.45
3:F:17:SER:CA	3:F:82:LEU:O	2.58	0.45
2:H:87:TYR:HD1	2:H:99:GLY:HA3	1.82	0.45
3:F:39:GLN:HG2	3:F:89:VAL:HG12	1.99	0.45
1:G:543:PHE:HD2	1:G:579:PRO:HD3	1.82	0.45
1:A:712:ILE:HG22	1:E:896:ILE:HD12	1.99	0.44
2:B:87:TYR:HD1	2:B:99:GLY:HA3	1.82	0.44
1:E:543:PHE:HD2	1:E:579:PRO:HD3	1.82	0.44
1:G:82:PRO:HG2	1:G:84:LEU:HD21	1.98	0.44
2:H:5:THR:HA	2:H:100:GLN:HG3	1.99	0.44
1:A:34:ARG:HH21	1:A:217:PRO:HB2	1.81	0.44
1:A:543:PHE:HD2	1:A:579:PRO:HD3	1.82	0.44
3:C:39:GLN:HG2	3:C:89:VAL:HG12	1.99	0.44
1:A:280:ASN:OD1	1:A:284:THR:N	2.49	0.44
1:A:570:ALA:HB1	1:E:963:VAL:HG11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:290:ASP:O	1:G:297:SER:HB3	2.18	0.44
1:G:598:ILE:HG23	1:G:664:ILE:HG21	1.99	0.44
2:D:79:GLU:H	2:D:82:ASP:HB2	1.83	0.44
2:D:27(A):SER:HA	2:D:69:THR:HG22	1.99	0.44
1:E:299:THR:HA	1:E:302:THR:HG22	2.00	0.44
1:E:598:ILE:HG23	1:E:664:ILE:HG21	1.99	0.44
3:F:10:GLU:CB	3:F:108:MET:O	2.53	0.43
3:F:20:VAL:HB	3:F:80:MET:SD	2.57	0.43
1:A:598:ILE:HG23	1:A:664:ILE:HG21	1.99	0.43
2:B:27(A):SER:HA	2:B:69:THR:HG22	1.99	0.43
1:E:715:PRO:HD3	1:G:894:LEU:HD13	1.99	0.43
3:I:96:ASN:HD21	3:I:102:LEU:HG	1.84	0.43
1:A:290:ASP:O	1:A:297:SER:HB3	2.18	0.43
1:A:914:ASN:ND2	1:G:1123:SER:OG	2.51	0.43
3:F:102:LEU:HB3	3:F:103:TRP:H	1.59	0.43
1:G:738:CYS:HB3	1:G:760:CYS:HB3	1.81	0.43
3:I:39:GLN:HG2	3:I:89:VAL:HG12	1.99	0.43
2:B:79:GLU:H	2:B:82:ASP:HB2	1.83	0.43
1:E:81:ASN:OD1	1:E:81:ASN:N	2.51	0.43
2:B:5:THR:HA	2:B:100:GLN:HG3	1.99	0.43
2:B:85:VAL:HA	2:B:102:THR:O	2.19	0.43
1:A:665:PRO:HA	1:A:671:CYS:SG	2.59	0.43
1:A:912:THR:O	1:A:915:VAL:HG22	2.19	0.43
1:A:963:VAL:HG11	1:G:570:ALA:HB1	2.01	0.43
2:D:5:THR:HA	2:D:100:GLN:HG3	1.99	0.43
1:G:81:ASN:N	1:G:81:ASN:OD1	2.51	0.43
2:D:85:VAL:HA	2:D:102:THR:O	2.18	0.43
1:E:665:PRO:HA	1:E:671:CYS:SG	2.59	0.43
2:H:79:GLU:H	2:H:82:ASP:HB2	1.83	0.43
2:H:85:VAL:HA	2:H:102:THR:O	2.18	0.43
1:E:290:ASP:O	1:E:297:SER:HB3	2.18	0.43
1:G:665:PRO:HA	1:G:671:CYS:SG	2.59	0.43
1:G:912:THR:O	1:G:915:VAL:HG22	2.19	0.43
3:I:97:CYS:HB2	3:I:100(B):CYS:HB2	1.30	0.43
1:A:28:TYR:HB3	1:A:61:ASN:OD1	2.19	0.42
1:A:81:ASN:OD1	1:A:81:ASN:N	2.51	0.42
1:A:299:THR:HA	1:A:302:THR:HG22	2.00	0.42
1:A:785:VAL:HG22	1:A:787:GLN:H	1.84	0.42
1:E:103:GLY:CA	1:E:241:LEU:HB2	2.47	0.42
1:E:611:LEU:HD22	1:E:666:ILE:HG23	2.00	0.42
1:E:912:THR:O	1:E:915:VAL:HG22	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:299:THR:HA	1:G:302:THR:HG22	2.00	0.42
1:A:103:GLY:CA	1:A:241:LEU:HB2	2.47	0.42
1:A:611:LEU:HD22	1:A:666:ILE:HG23	2.00	0.42
1:E:28:TYR:HB3	1:E:61:ASN:OD1	2.19	0.42
1:E:785:VAL:HG22	1:E:787:GLN:H	1.84	0.42
1:A:740:MET:HB2	1:G:319:ARG:NH2	2.35	0.42
3:F:97:CYS:HB2	3:F:100(C):TYR:H	1.85	0.42
3:I:97:CYS:HB2	3:I:100(C):TYR:H	1.85	0.42
1:A:676:THR:HA	1:A:690:GLN:HA	2.01	0.42
1:A:906:PHE:HD2	1:A:916:LEU:HB2	1.85	0.42
2:B:54:ARG:NH1	2:B:63:SER:HA	2.34	0.42
1:G:28:TYR:HB3	1:G:61:ASN:OD1	2.19	0.42
1:G:785:VAL:HG22	1:G:787:GLN:H	1.84	0.42
3:F:19:LYS:HZ1	3:F:79:TYR:HB3	1.84	0.42
3:I:52:VAL:O	3:I:55:GLY:CA	2.68	0.42
1:G:676:THR:HA	1:G:690:GLN:HA	2.01	0.42
1:G:906:PHE:HD2	1:G:916:LEU:HB2	1.85	0.42
3:C:96:ASN:HD21	3:C:102:LEU:HG	1.84	0.41
3:F:97:CYS:HB2	3:F:100(B):CYS:HB2	1.30	0.41
2:D:54:ARG:NH1	2:D:63:SER:HA	2.34	0.41
1:G:194:PHE:HD1	1:G:203:ILE:HG12	1.85	0.41
3:C:97:CYS:HB2	3:C:100(B):CYS:HB2	1.30	0.41
2:D:63:SER:O	2:D:73:LEU:HD12	2.21	0.41
1:E:194:PHE:HD1	1:E:203:ILE:HG12	1.85	0.41
1:E:676:THR:HA	1:E:690:GLN:HA	2.01	0.41
1:G:103:GLY:CA	1:G:241:LEU:HB2	2.47	0.41
1:G:193:VAL:HG23	1:G:223:LEU:HD22	2.02	0.41
1:G:611:LEU:HD22	1:G:666:ILE:HG23	2.00	0.41
2:B:63:SER:O	2:B:73:LEU:HD12	2.21	0.41
1:E:906:PHE:HD2	1:E:916:LEU:HB2	1.85	0.41
1:E:1107:ARG:HG2	1:G:904:TYR:OH	2.20	0.41
3:F:52:VAL:O	3:F:55:GLY:CA	2.68	0.41
2:H:63:SER:O	2:H:73:LEU:HD12	2.21	0.41
1:A:917:TYR:HB3	1:G:1129:VAL:HG23	2.01	0.41
3:C:97:CYS:HB2	3:C:100(C):TYR:H	1.85	0.41
1:A:193:VAL:HG23	1:A:223:LEU:HD22	2.02	0.41
1:A:206:LYS:HB2	1:A:223:LEU:HA	2.03	0.41
3:C:52:VAL:O	3:C:55:GLY:CA	2.68	0.41
1:E:206:LYS:HB2	1:E:223:LEU:HA	2.03	0.41
1:E:980:ILE:HD13	1:E:992:GLN:HB3	2.03	0.41
1:A:738:CYS:HB3	1:A:760:CYS:HB3	1.81	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:980:ILE:HD13	1:A:992:GLN:HB3	2.03	0.41
1:E:738:CYS:HB3	1:E:760:CYS:HB3	1.81	0.41
1:A:1080:ALA:O	1:A:1132:ILE:HG13	2.21	0.40
1:E:1123:SER:OG	1:G:914:ASN:ND2	2.54	0.40
1:G:40:ASP:OD1	1:G:41:LYS:N	2.52	0.40
1:G:206:LYS:HB2	1:G:223:LEU:HA	2.03	0.40
3:C:35:GLN:OE1	2:D:98:PHE:CZ	2.75	0.40
1:A:194:PHE:HD1	1:A:203:ILE:HG12	1.85	0.40
3:C:32:SER:OG	3:C:95:PRO:HD2	2.22	0.40
1:E:722:VAL:HG22	1:E:1065:VAL:HG22	2.03	0.40
3:C:19:LYS:HZ1	3:C:79:TYR:HB3	1.87	0.40
2:H:54:ARG:NH1	2:H:63:SER:HA	2.34	0.40
1:A:310:LYS:NZ	1:A:663:ASP:OD2	2.37	0.40
1:E:669:GLY:HA2	1:E:697:MET:HE2	2.03	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	982/1288 (76%)	955 (97%)	27 (3%)	0	100	100
1	E	982/1288 (76%)	955 (97%)	27 (3%)	0	100	100
1	G	982/1288 (76%)	955 (97%)	27 (3%)	0	100	100
2	B	106/215 (49%)	100 (94%)	6 (6%)	0	100	100
2	D	106/215 (49%)	100 (94%)	6 (6%)	0	100	100
2	H	106/215 (49%)	100 (94%)	6 (6%)	0	100	100
3	C	121/231 (52%)	111 (92%)	10 (8%)	0	100	100
3	F	121/231 (52%)	111 (92%)	10 (8%)	0	100	100
3	I	121/231 (52%)	111 (92%)	10 (8%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	3627/5202 (70%)	3498 (96%)	129 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	631/1116 (56%)	631 (100%)	0	100	100
1	E	631/1116 (56%)	631 (100%)	0	100	100
1	G	631/1116 (56%)	631 (100%)	0	100	100
2	B	90/186 (48%)	89 (99%)	1 (1%)	73	85
2	D	90/186 (48%)	89 (99%)	1 (1%)	73	85
2	H	90/186 (48%)	89 (99%)	1 (1%)	73	85
3	C	102/196 (52%)	101 (99%)	1 (1%)	76	86
3	F	102/196 (52%)	101 (99%)	1 (1%)	76	86
3	I	102/196 (52%)	101 (99%)	1 (1%)	76	86
All	All	2469/4494 (55%)	2463 (100%)	6 (0%)	93	97

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

Mol	Chain	Res	Type
2	B	24	ARG
3	C	12	LYS
2	D	24	ARG
3	F	12	LYS
2	H	24	ARG
3	I	12	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	913	GLN
3	C	35	GLN
3	C	96	ASN
1	E	913	GLN
3	F	96	ASN
1	G	913	GLN
3	I	96	ASN

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 monosaccharides 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.

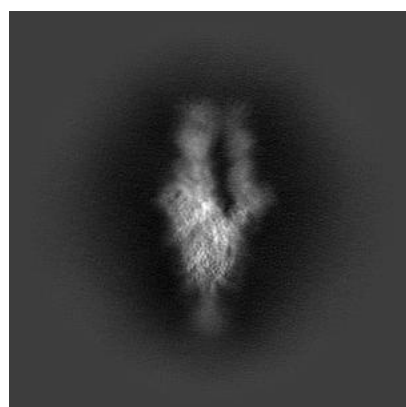
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12645. These allow visual inspection of the internal detail of the map and identification of artifacts.

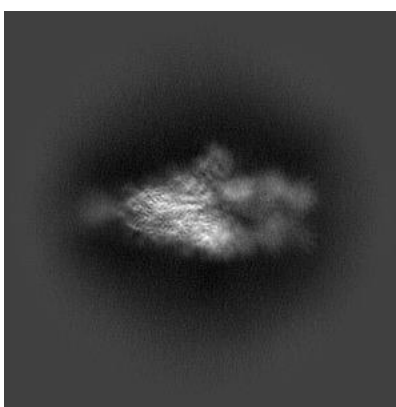
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

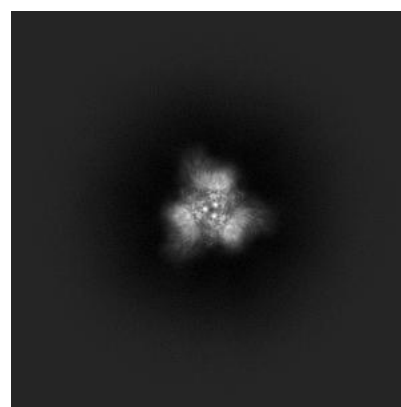
6.1.1 Primary map



X



Y



Z

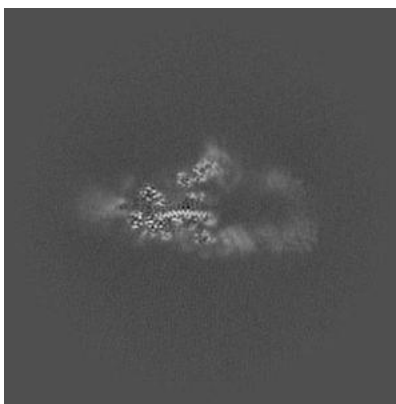
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

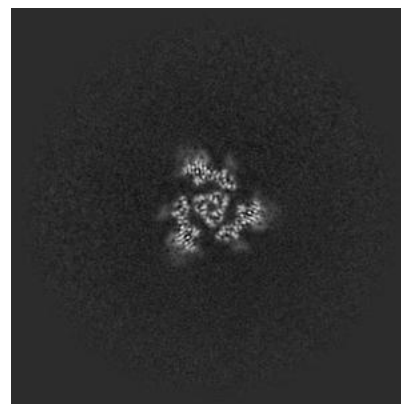
6.2.1 Primary map



X Index: 150



Y Index: 150

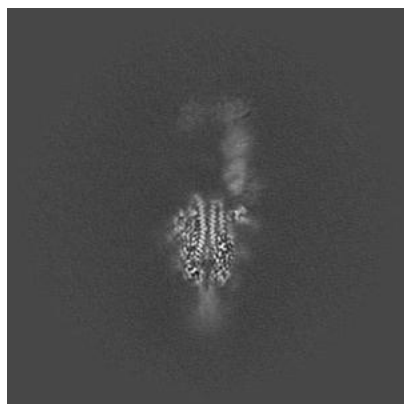


Z Index: 150

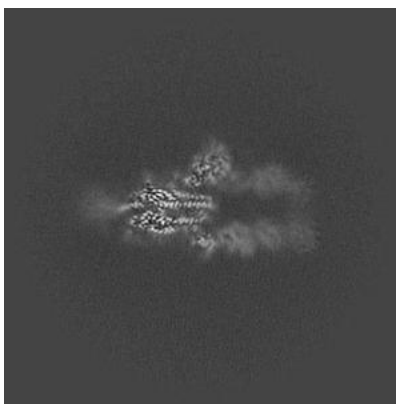
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

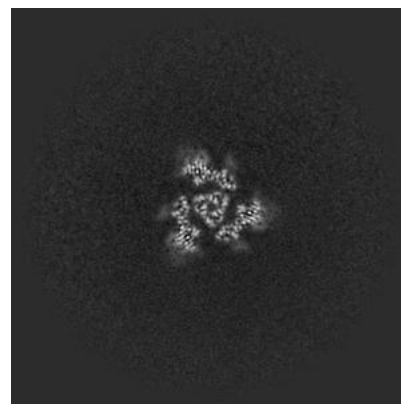
6.3.1 Primary map



X Index: 152



Y Index: 145



Z Index: 150

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

6.4 Orthogonal surface views [i](#)

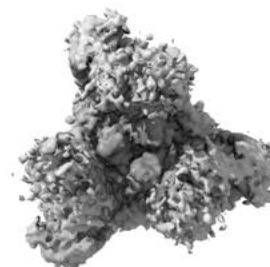
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.3. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

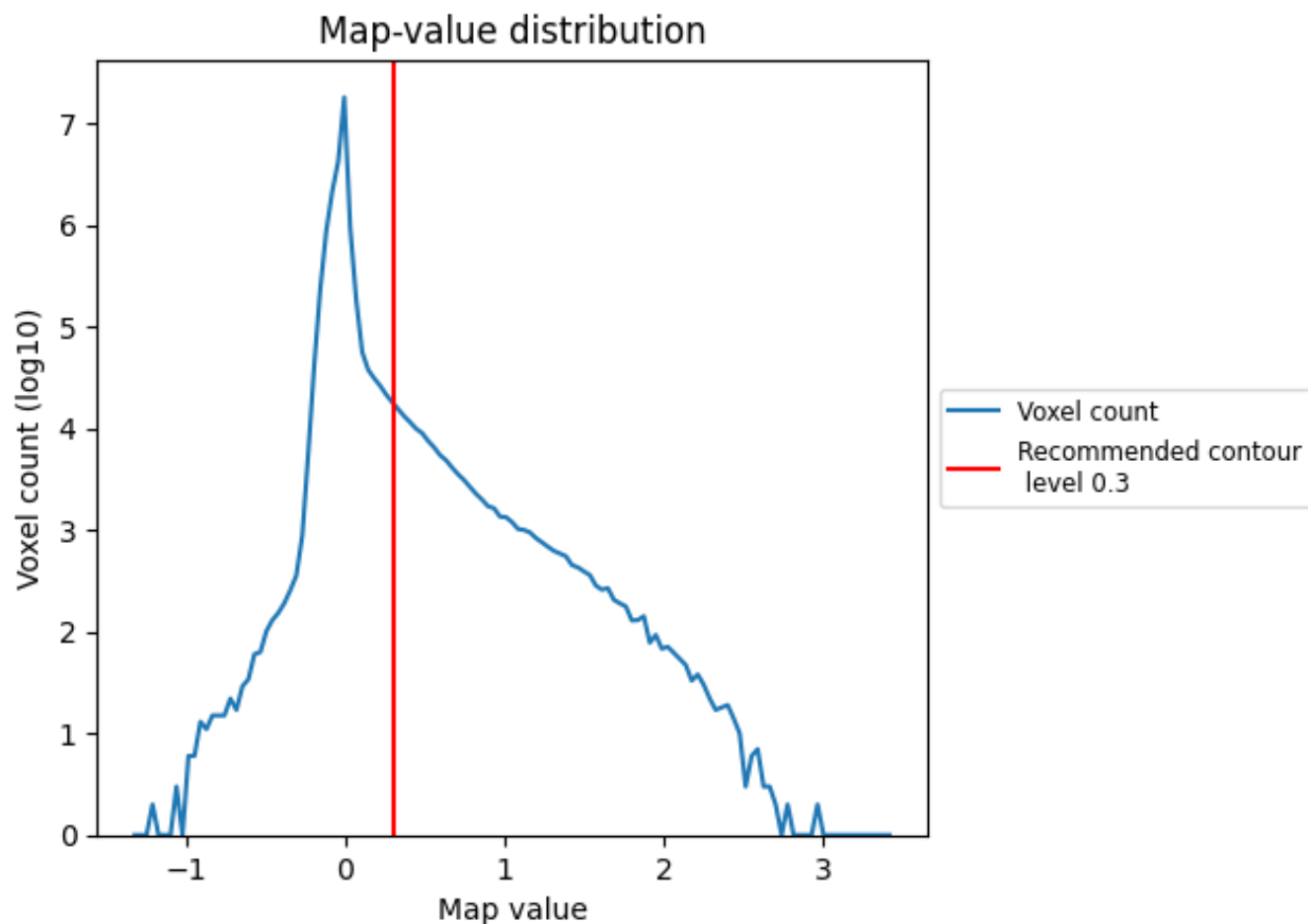
6.5 Mask visualisation

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

7 Map analysis [i](#)

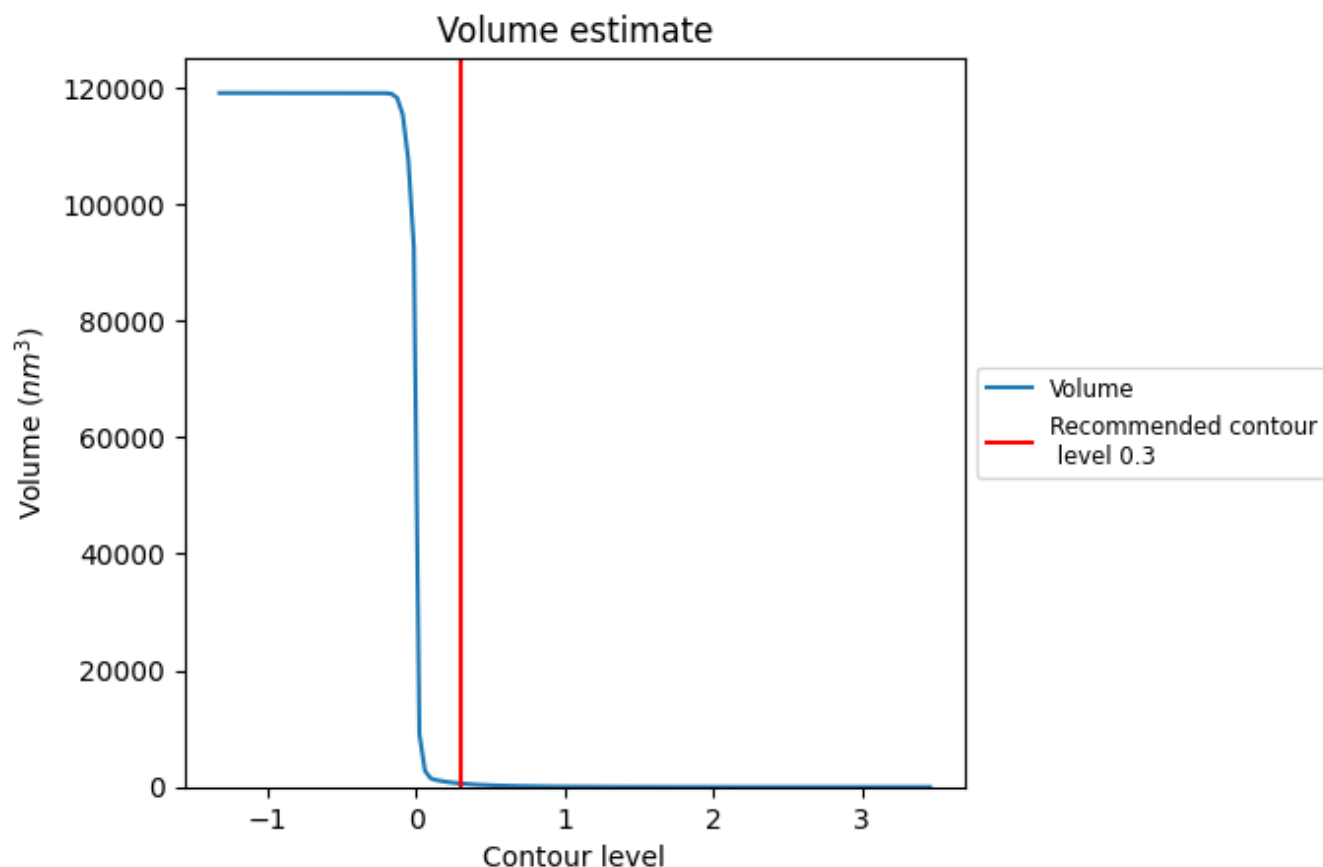
This section contains the results of statistical analysis of the map.

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.

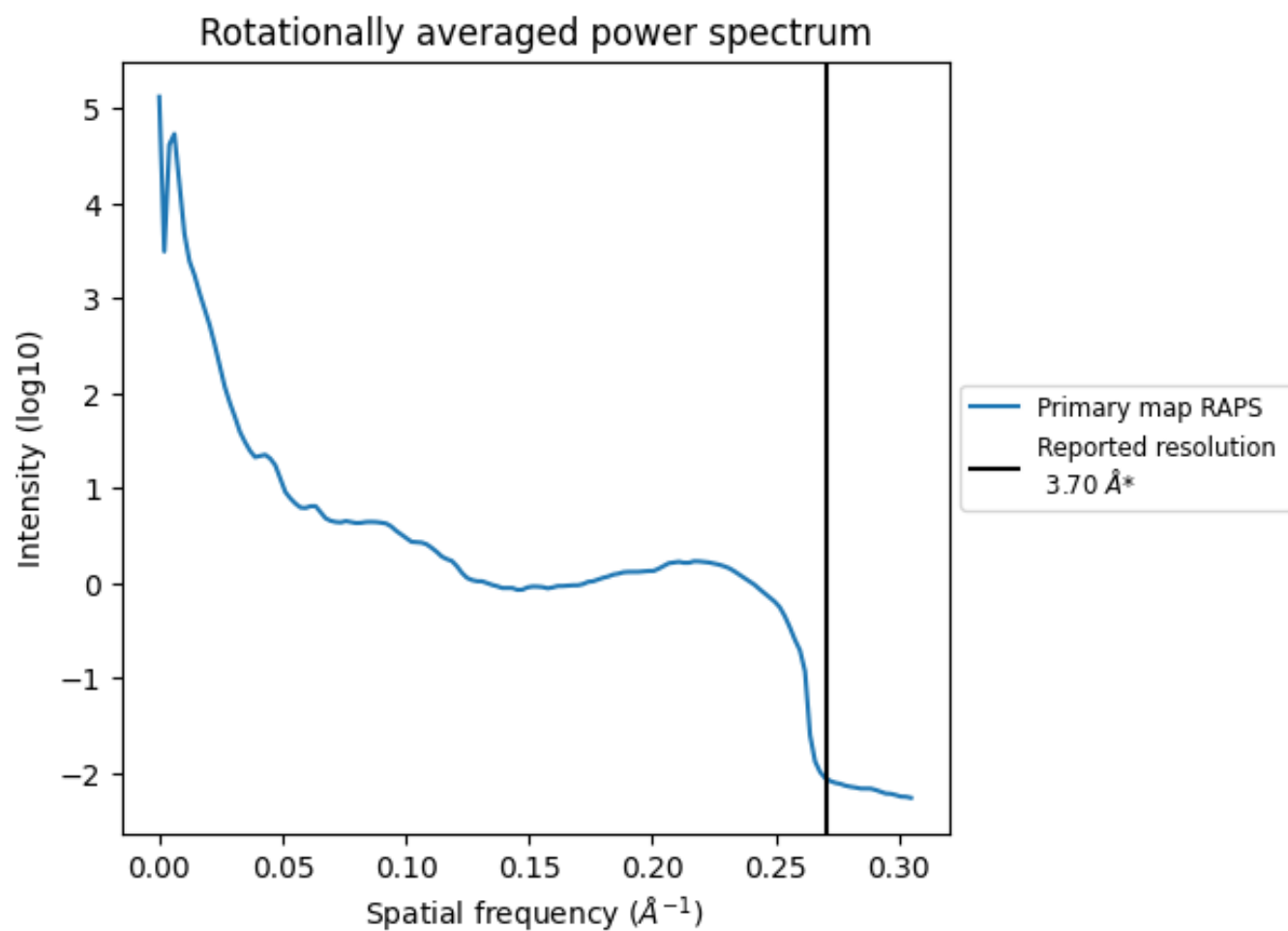
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 601 nm^3 ; this corresponds to an approximate mass of 543 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 ⓘ

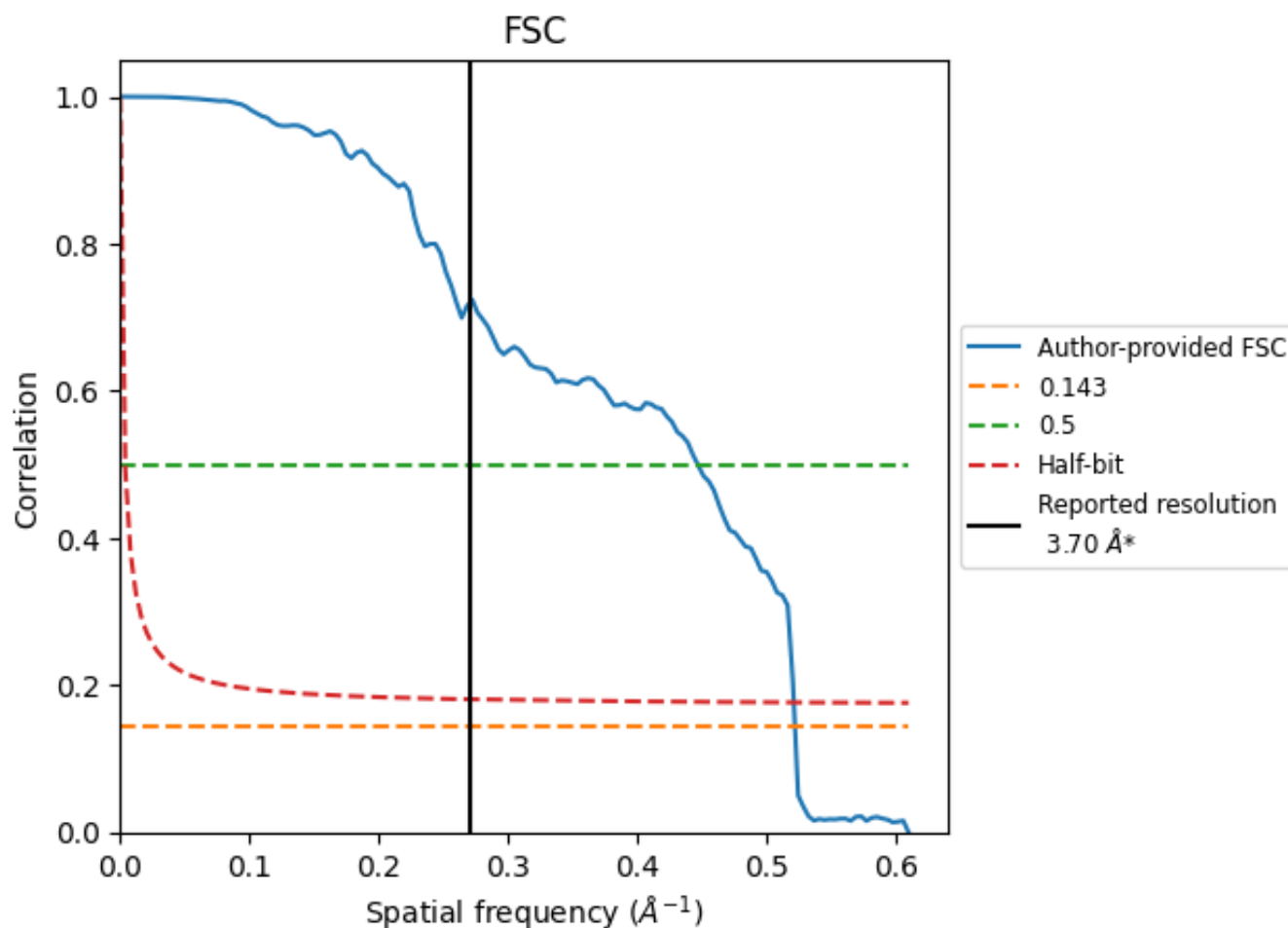


*Reported resolution corresponds to spatial frequency of 0.270 Å⁻¹

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 0.270 \AA^{-1}

8.2 Resolution estimates [i](#)

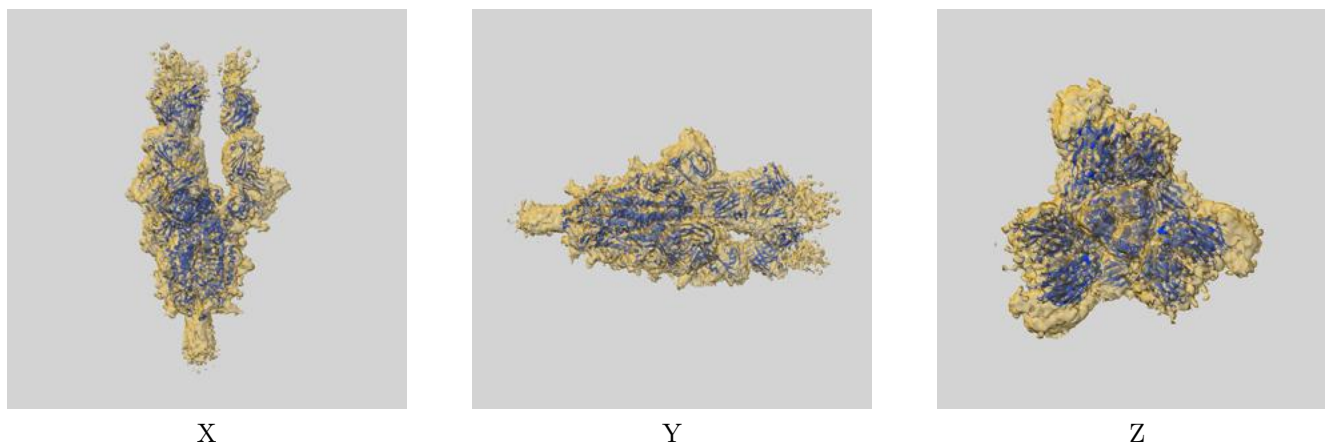
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	1.92	2.24	1.92
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from author-provided FSC intersecting FSC 0.143 CUT-OFF 1.92 differs from the reported value 3.7 by more than 10 %

9 Map-model fit [i](#)

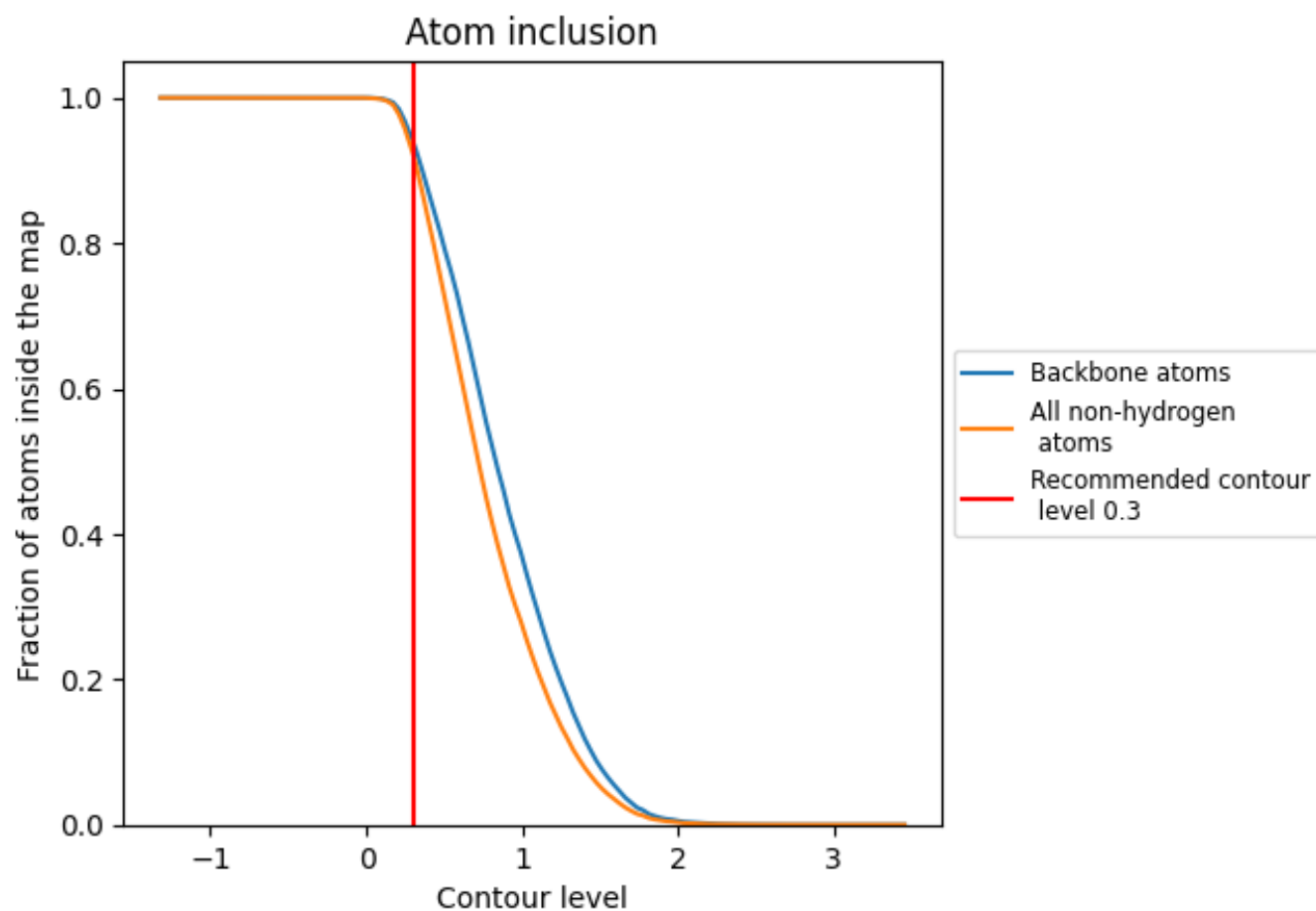
This section contains information regarding the fit between EMDB map EMD-12645 and PDB model 7NY5. Per-residue inclusion information can be found in section [3](#) on page [12](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.3 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 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 92% of all non-hydrogen atoms, are inside the map.