



Full wwPDB EM Validation Report ⓘ

Mar 25, 2026 – 07:47 AM UTC

PDB ID : 8H3M / pdb_00008h3m
EMDB ID : EMD-34469
Title : Conformation 1 of SARS-CoV-2 Omicron BA.1 Variant Spike protein complexed with MO1 Fab
Authors : Ishimaru, H.; Nishimura, M.; Sutandhio, S.; Shigematsu, H.; Kato, K.; Hasegawa, N.; Mori, Y.
Deposited on : 2022-10-09
Resolution : 2.48 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

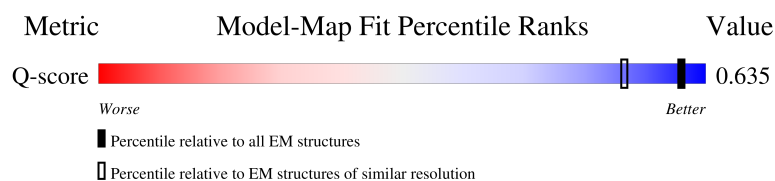
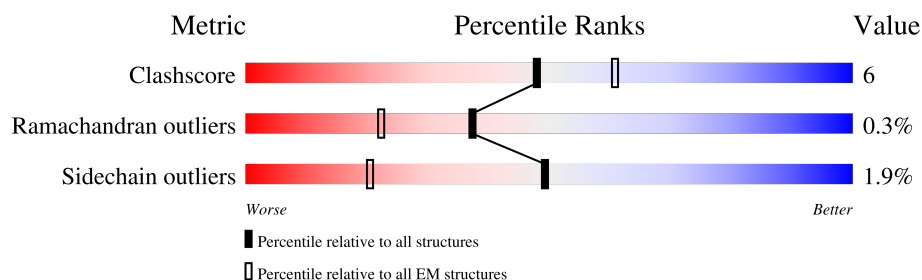
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

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

The reported resolution of this entry is 2.48 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	6178 (1.98 - 2.98)

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	1249	
1	B	1249	
1	C	1249	
2	D	449	

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Mol	Chain	Length	Quality of chain
2	H	449	 94%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20936 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	910	Total	C	N	O	S	0	0
			7039	4521	1173	1314	31		
1	B	763	Total	C	N	O	S	0	0
			5883	3777	979	1102	25		
1	C	930	Total	C	N	O	S	0	0
			7213	4633	1208	1341	31		

There are 279 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	VAL	ALA	variant	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	95	ILE	THR	variant	UNP P0DTC2
A	?	-	GLY	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	145	ASP	TYR	variant	UNP P0DTC2
A	209B	ILE	-	insertion	UNP P0DTC2
A	209C	VAL	-	insertion	UNP P0DTC2
A	209D	ARG	ASN	variant	UNP P0DTC2
A	209E	GLU	LEU	variant	UNP P0DTC2
A	209F	PRO	VAL	variant	UNP P0DTC2
A	209G	GLU	ARG	variant	UNP P0DTC2
A	339	ASP	GLY	variant	UNP P0DTC2
A	371	LEU	SER	variant	UNP P0DTC2
A	373	PRO	SER	variant	UNP P0DTC2
A	375	PHE	SER	variant	UNP P0DTC2
A	417	ASN	LYS	variant	UNP P0DTC2
A	440	LYS	ASN	variant	UNP P0DTC2
A	446	SER	GLY	variant	UNP P0DTC2
A	477	ASN	SER	variant	UNP P0DTC2
A	478	LYS	THR	variant	UNP P0DTC2
A	484	ALA	GLU	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	493	ARG	GLN	variant	UNP P0DTC2
A	496	SER	GLY	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	547	LYS	THR	variant	UNP P0DTC2
A	614	GLY	ASP	variant	UNP P0DTC2
A	655	TYR	HIS	variant	UNP P0DTC2
A	682	LYS	ASN	variant	UNP P0DTC2
A	?	-	PRO	deletion	UNP P0DTC2
A	?	-	ARG	deletion	UNP P0DTC2
A	?	-	ARG	deletion	UNP P0DTC2
A	684	HIS	ALA	variant	UNP P0DTC2
A	685	ALA	ARG	variant	UNP P0DTC2
A	764	LYS	ASN	variant	UNP P0DTC2
A	796	TYR	ASP	variant	UNP P0DTC2
A	817	PRO	PHE	variant	UNP P0DTC2
A	856	LYS	ASN	variant	UNP P0DTC2
A	892	PRO	ALA	variant	UNP P0DTC2
A	899	PRO	ALA	variant	UNP P0DTC2
A	942	PRO	ALA	variant	UNP P0DTC2
A	954	HIS	GLN	variant	UNP P0DTC2
A	969	LYS	ASN	variant	UNP P0DTC2
A	981	PHE	LEU	variant	UNP P0DTC2
A	986	PRO	LYS	variant	UNP P0DTC2
A	987	PRO	VAL	variant	UNP P0DTC2
A	1213	PRO	-	expression tag	UNP P0DTC2
A	1214	LEU	-	expression tag	UNP P0DTC2
A	1215	GLU	-	expression tag	UNP P0DTC2
A	1216	VAL	-	expression tag	UNP P0DTC2
A	1217	LEU	-	expression tag	UNP P0DTC2
A	1218	PHE	-	expression tag	UNP P0DTC2
A	1219	GLN	-	expression tag	UNP P0DTC2
A	1220	GLY	-	expression tag	UNP P0DTC2
A	1221	PRO	-	expression tag	UNP P0DTC2
A	1222	GLY	-	expression tag	UNP P0DTC2
A	1223	TYR	-	expression tag	UNP P0DTC2
A	1224	ILE	-	expression tag	UNP P0DTC2
A	1225	PRO	-	expression tag	UNP P0DTC2
A	1226	GLU	-	expression tag	UNP P0DTC2
A	1227	ALA	-	expression tag	UNP P0DTC2
A	1228	PRO	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1229	ARG	-	expression tag	UNP P0DTC2
A	1230	ASP	-	expression tag	UNP P0DTC2
A	1231	GLY	-	expression tag	UNP P0DTC2
A	1232	GLN	-	expression tag	UNP P0DTC2
A	1233	ALA	-	expression tag	UNP P0DTC2
A	1234	TYR	-	expression tag	UNP P0DTC2
A	1235	VAL	-	expression tag	UNP P0DTC2
A	1236	ARG	-	expression tag	UNP P0DTC2
A	1237	LYS	-	expression tag	UNP P0DTC2
A	1238	ASP	-	expression tag	UNP P0DTC2
A	1239	GLY	-	expression tag	UNP P0DTC2
A	1240	GLU	-	expression tag	UNP P0DTC2
A	1241	TRP	-	expression tag	UNP P0DTC2
A	1242	VAL	-	expression tag	UNP P0DTC2
A	1243	PHE	-	expression tag	UNP P0DTC2
A	1244	LEU	-	expression tag	UNP P0DTC2
A	1245	SER	-	expression tag	UNP P0DTC2
A	1246	THR	-	expression tag	UNP P0DTC2
A	1247	PHE	-	expression tag	UNP P0DTC2
A	1248	LEU	-	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
B	69	VAL	ALA	variant	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	95	ILE	THR	variant	UNP P0DTC2
B	?	-	GLY	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	145	ASP	TYR	variant	UNP P0DTC2
B	207D	ILE	-	insertion	UNP P0DTC2
B	207E	VAL	-	insertion	UNP P0DTC2
B	207F	ARG	ASN	variant	UNP P0DTC2
B	207G	GLU	LEU	variant	UNP P0DTC2
B	207H	PRO	VAL	variant	UNP P0DTC2
B	207I	GLU	ARG	variant	UNP P0DTC2
B	339	ASP	GLY	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
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	682	LYS	ASN	variant	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	?	-	ARG	deletion	UNP P0DTC2
B	?	-	ARG	deletion	UNP P0DTC2
B	684	HIS	ALA	variant	UNP P0DTC2
B	685	ALA	ARG	variant	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	817	PRO	PHE	variant	UNP P0DTC2
B	856	LYS	ASN	variant	UNP P0DTC2
B	892	PRO	ALA	variant	UNP P0DTC2
B	899	PRO	ALA	variant	UNP P0DTC2
B	942	PRO	ALA	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	1213	PRO	-	expression tag	UNP P0DTC2
B	1214	LEU	-	expression tag	UNP P0DTC2
B	1215	GLU	-	expression tag	UNP P0DTC2
B	1216	VAL	-	expression tag	UNP P0DTC2
B	1217	LEU	-	expression tag	UNP P0DTC2
B	1218	PHE	-	expression tag	UNP P0DTC2
B	1219	GLN	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1220	GLY	-	expression tag	UNP P0DTC2
B	1221	PRO	-	expression tag	UNP P0DTC2
B	1222	GLY	-	expression tag	UNP P0DTC2
B	1223	TYR	-	expression tag	UNP P0DTC2
B	1224	ILE	-	expression tag	UNP P0DTC2
B	1225	PRO	-	expression tag	UNP P0DTC2
B	1226	GLU	-	expression tag	UNP P0DTC2
B	1227	ALA	-	expression tag	UNP P0DTC2
B	1228	PRO	-	expression tag	UNP P0DTC2
B	1229	ARG	-	expression tag	UNP P0DTC2
B	1230	ASP	-	expression tag	UNP P0DTC2
B	1231	GLY	-	expression tag	UNP P0DTC2
B	1232	GLN	-	expression tag	UNP P0DTC2
B	1233	ALA	-	expression tag	UNP P0DTC2
B	1234	TYR	-	expression tag	UNP P0DTC2
B	1235	VAL	-	expression tag	UNP P0DTC2
B	1236	ARG	-	expression tag	UNP P0DTC2
B	1237	LYS	-	expression tag	UNP P0DTC2
B	1238	ASP	-	expression tag	UNP P0DTC2
B	1239	GLY	-	expression tag	UNP P0DTC2
B	1240	GLU	-	expression tag	UNP P0DTC2
B	1241	TRP	-	expression tag	UNP P0DTC2
B	1242	VAL	-	expression tag	UNP P0DTC2
B	1243	PHE	-	expression tag	UNP P0DTC2
B	1244	LEU	-	expression tag	UNP P0DTC2
B	1245	SER	-	expression tag	UNP P0DTC2
B	1246	THR	-	expression tag	UNP P0DTC2
B	1247	PHE	-	expression tag	UNP P0DTC2
B	1248	LEU	-	expression tag	UNP P0DTC2
B	1249	GLY	-	expression tag	UNP P0DTC2
B	1250	HIS	-	expression tag	UNP P0DTC2
B	1251	HIS	-	expression tag	UNP P0DTC2
B	1252	HIS	-	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
C	67	VAL	ALA	variant	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	95	ILE	THR	variant	UNP P0DTC2
C	?	-	GLY	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	TYR	deletion	UNP P0DTC2
C	145	ASP	TYR	variant	UNP P0DTC2
C	209B	ILE	-	insertion	UNP P0DTC2
C	209C	VAL	-	insertion	UNP P0DTC2
C	209D	ARG	ASN	variant	UNP P0DTC2
C	209E	GLU	LEU	variant	UNP P0DTC2
C	209F	PRO	VAL	variant	UNP P0DTC2
C	209G	GLU	ARG	variant	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	682	LYS	ASN	variant	UNP P0DTC2
C	?	-	PRO	deletion	UNP P0DTC2
C	?	-	ARG	deletion	UNP P0DTC2
C	?	-	ARG	deletion	UNP P0DTC2
C	684	HIS	ALA	variant	UNP P0DTC2
C	685	ALA	ARG	variant	UNP P0DTC2
C	764	LYS	ASN	variant	UNP P0DTC2
C	796	TYR	ASP	variant	UNP P0DTC2
C	817	PRO	PHE	variant	UNP P0DTC2
C	856	LYS	ASN	variant	UNP P0DTC2
C	892	PRO	ALA	variant	UNP P0DTC2
C	899	PRO	ALA	variant	UNP P0DTC2
C	942	PRO	ALA	variant	UNP P0DTC2
C	954	HIS	GLN	variant	UNP P0DTC2
C	969	LYS	ASN	variant	UNP P0DTC2
C	981	PHE	LEU	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	986	PRO	LYS	variant	UNP P0DTC2
C	987	PRO	VAL	variant	UNP P0DTC2
C	1213	PRO	-	expression tag	UNP P0DTC2
C	1214	LEU	-	expression tag	UNP P0DTC2
C	1215	GLU	-	expression tag	UNP P0DTC2
C	1216	VAL	-	expression tag	UNP P0DTC2
C	1217	LEU	-	expression tag	UNP P0DTC2
C	1218	PHE	-	expression tag	UNP P0DTC2
C	1219	GLN	-	expression tag	UNP P0DTC2
C	1220	GLY	-	expression tag	UNP P0DTC2
C	1221	PRO	-	expression tag	UNP P0DTC2
C	1222	GLY	-	expression tag	UNP P0DTC2
C	1223	TYR	-	expression tag	UNP P0DTC2
C	1224	ILE	-	expression tag	UNP P0DTC2
C	1225	PRO	-	expression tag	UNP P0DTC2
C	1226	GLU	-	expression tag	UNP P0DTC2
C	1227	ALA	-	expression tag	UNP P0DTC2
C	1228	PRO	-	expression tag	UNP P0DTC2
C	1229	ARG	-	expression tag	UNP P0DTC2
C	1230	ASP	-	expression tag	UNP P0DTC2
C	1231	GLY	-	expression tag	UNP P0DTC2
C	1232	GLN	-	expression tag	UNP P0DTC2
C	1233	ALA	-	expression tag	UNP P0DTC2
C	1234	TYR	-	expression tag	UNP P0DTC2
C	1235	VAL	-	expression tag	UNP P0DTC2
C	1236	ARG	-	expression tag	UNP P0DTC2
C	1237	LYS	-	expression tag	UNP P0DTC2
C	1238	ASP	-	expression tag	UNP P0DTC2
C	1239	GLY	-	expression tag	UNP P0DTC2
C	1240	GLU	-	expression tag	UNP P0DTC2
C	1241	TRP	-	expression tag	UNP P0DTC2
C	1242	VAL	-	expression tag	UNP P0DTC2
C	1243	PHE	-	expression tag	UNP P0DTC2
C	1244	LEU	-	expression tag	UNP P0DTC2
C	1245	SER	-	expression tag	UNP P0DTC2
C	1246	THR	-	expression tag	UNP P0DTC2
C	1247	PHE	-	expression tag	UNP P0DTC2
C	1248	LEU	-	expression tag	UNP P0DTC2
C	1249	GLY	-	expression tag	UNP P0DTC2
C	1250	HIS	-	expression tag	UNP P0DTC2
C	1251	HIS	-	expression tag	UNP P0DTC2
C	1252	HIS	-	expression tag	UNP P0DTC2

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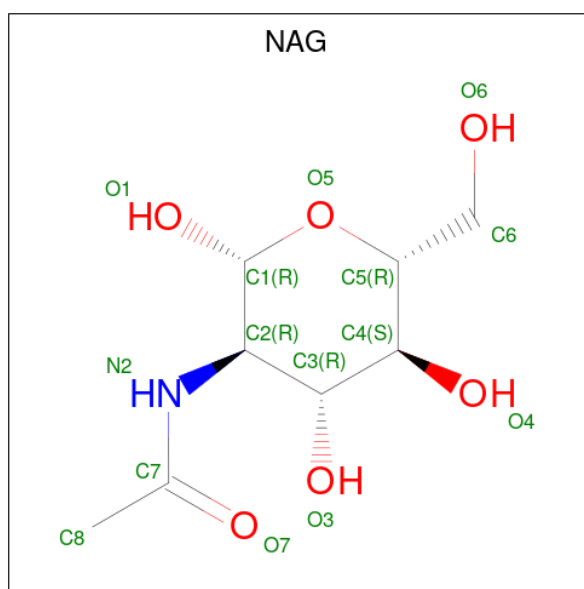
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Chain	Residue	Modelled	Actual	Comment	Reference
C	1253	HIS	-	expression tag	UNP P0DTC2
C	1254	HIS	-	expression tag	UNP P0DTC2
C	1255	HIS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called MO1 heavy chain.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	D	30	Total	C	N	O	0	0
			255	164	38	53		
2	H	29	Total	C	N	O	S	0
			238	149	38	50	1	0

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

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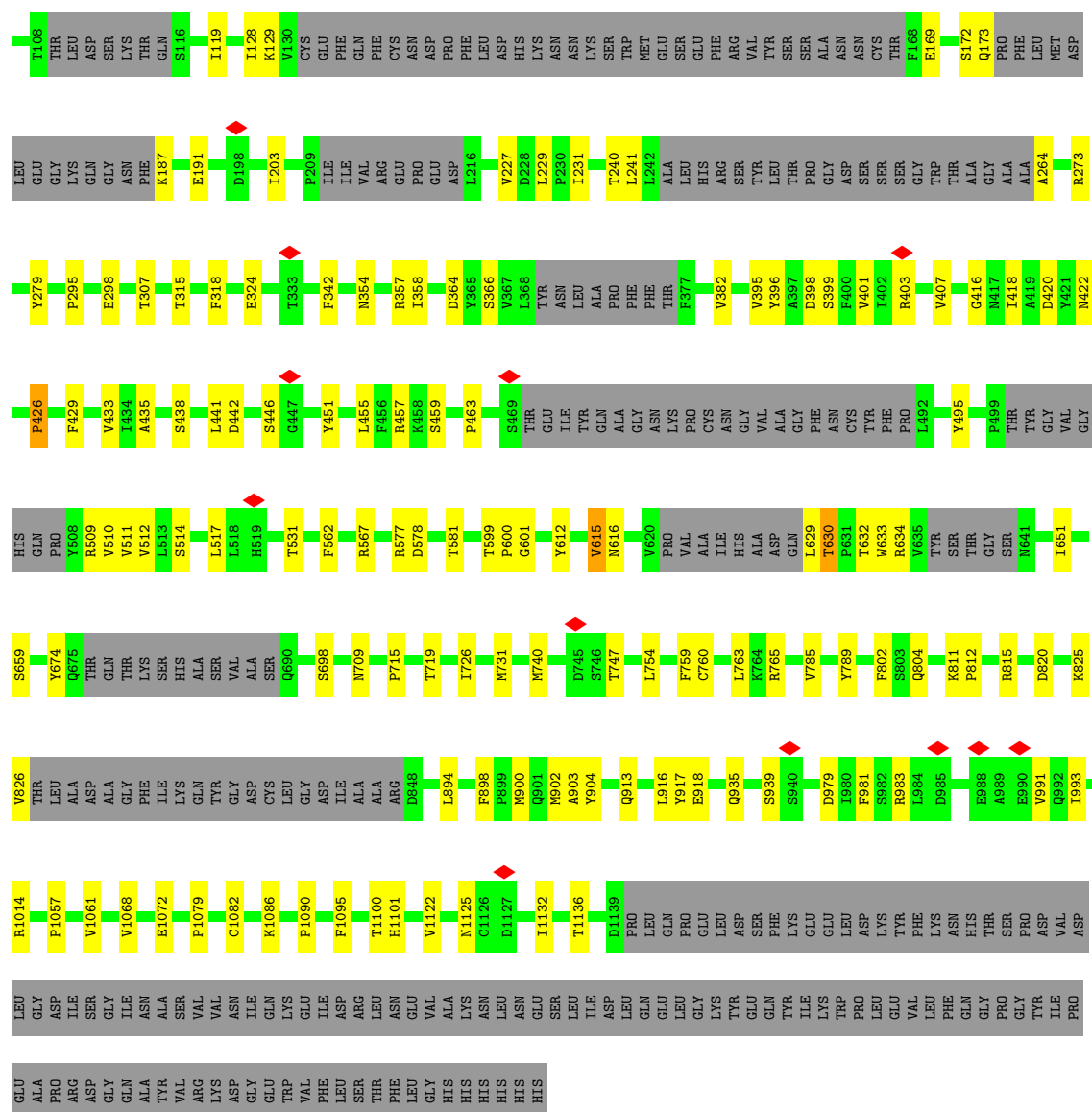
Mol	Chain	Residues	Atoms				AltConf
3	A	1	Total	C	N	O	0
			14	8	1	5	
3	A	1	Total	C	N	O	0
			14	8	1	5	
3	A	1	Total	C	N	O	0
			14	8	1	5	
3	A	1	Total	C	N	O	0
			14	8	1	5	
3	B	1	Total	C	N	O	0
			14	8	1	5	
3	B	1	Total	C	N	O	0
			14	8	1	5	
3	B	1	Total	C	N	O	0
			14	8	1	5	
3	B	1	Total	C	N	O	0
			14	8	1	5	
3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

Category	Percentage
Very bad	51%
Bad	10%
Good	39%



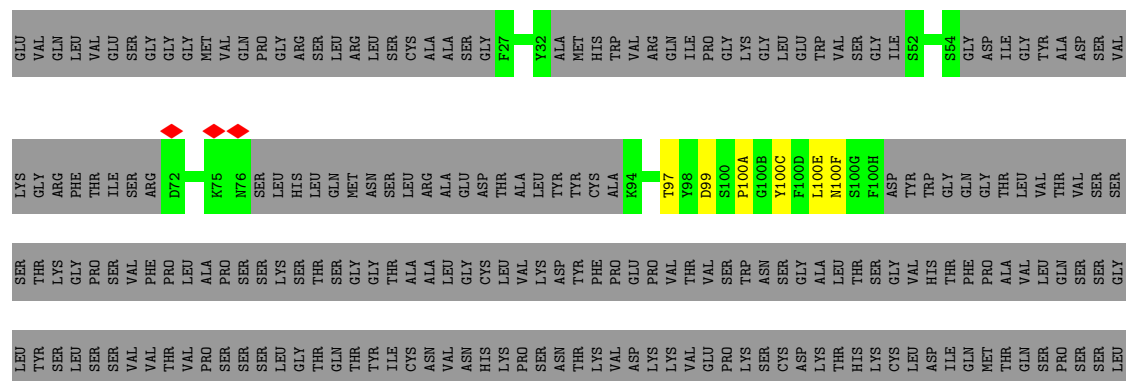
Response	Percentage
Yes	63%
No	12%
Don't know	26%





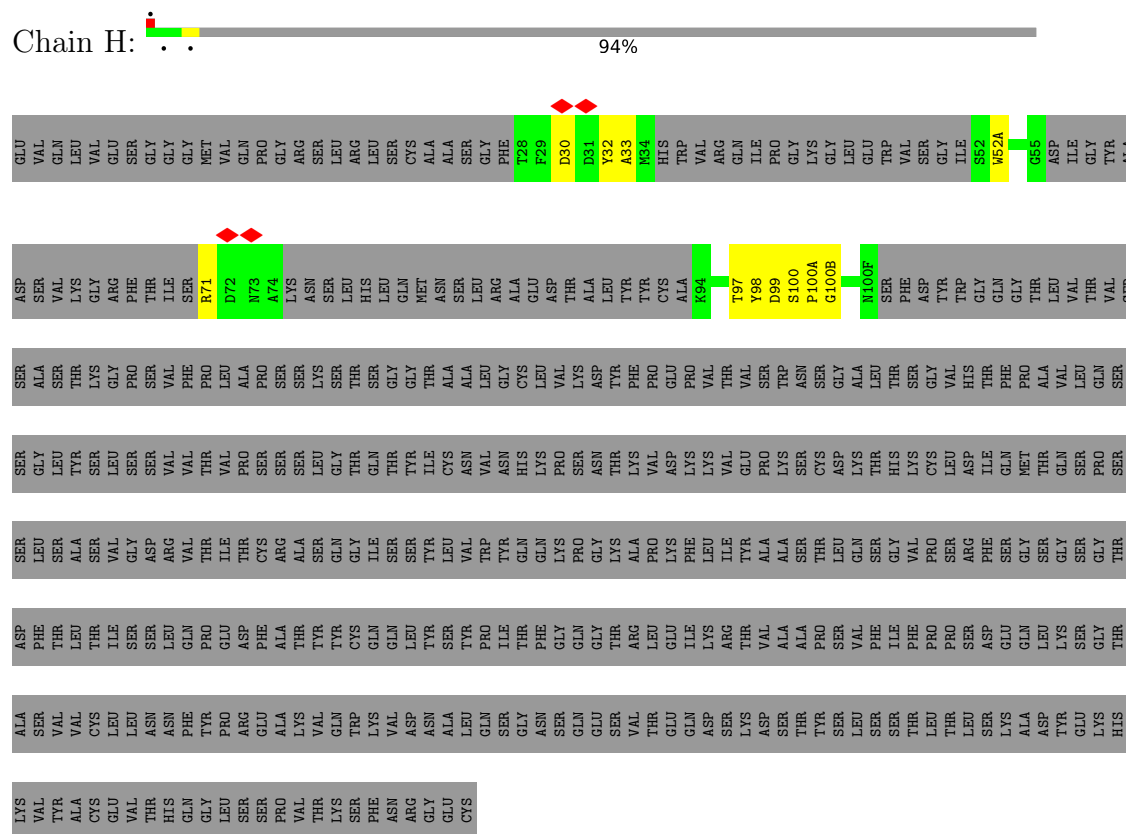
• Molecule 2: MO1 heavy chain

Chain D: 5% . 93%



Tyr	Val	Thr	Ser
Ala	Val	Leu	Ala
Cys	Cys	Thr	Ser
Glu	Leu	Ile	Val
Val	Leu	Ser	Gly
Thr	Asn	Ser	Asp
His	Asn	Leu	Arg
Gln	Phe	Gln	Val
Gly	Tyr	Pro	Thr
Leu	Pro	Glu	Ile
Ser	Arg	Asp	Thr
Ser	Glu	Phe	Cys
Pro	Ala	Ala	Arg
Val	Lys	Thr	Ala
Thr	Val	Tyr	Ser
Lys	Gln	Tyr	Ser
Ser	Trp	Cys	Gly
Phe	Lys	Gln	Ile
Asn	Val	Gln	Ser
Arg	Asp	Leu	Ser
Gly	Asn	Tyr	Tyr
Glu	Ala	Ser	Leu
Cys	Leu	Tyr	Val
	Gln	Pro	Trp
	Ser	Ile	Tyr
	Gly	Thr	Gln
	Asn	Phe	Gln
	Ser	Gly	Lys
	Gln	Gln	Pro
	Glu	Gly	Gly
	Ser	Thr	Lys
	Val	Arg	Ala
	Thr	Leu	Pro
	Glu	Ile	Lys
	Asp	Ser	Phe
	Val	Thr	Leu
	Gln	Arg	Thr
	Ala	Ala	Ser
	Asn	Pro	Thr
	Ser	Val	Leu
	Thr	Phe	Gln
	Ser	Ile	Gly
	Thr	Thr	Val
	Leu	Pro	Pro
	Thr	Pro	Ser
	Ser	Asp	Arg
	Leu	Ser	Phe
	Lys	Ser	Ser
	Ala	Glu	Gly
	Asp	Leu	Ser
	Tyr	Lys	Gly
	Glu	Ser	Gly
	Lys	Thr	Thr
	His	Thr	Asp
	Val	Ser	Phe

- Molecule 2: MO1 heavy chain



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	533315	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	JEOL CRYO ARM 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	1600	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.062	Depositor
Minimum map value	-0.039	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.005	Depositor
Map size ($\text{\AA}$)	385.024, 385.024, 385.024	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.752, 0.752, 0.752	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.16	0/7199	0.36	1/9807 (0.0%)
1	B	0.16	0/6009	0.41	0/8172
1	C	0.14	0/7372	0.35	2/10028 (0.0%)
2	D	0.14	0/261	0.50	0/348
2	H	0.16	0/242	0.46	0/322
All	All	0.15	0/21083	0.37	3/28677 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	426	PRO	N-CD-CG	-6.73	93.11	103.20
1	A	521	PRO	CA-N-CD	-6.58	102.79	112.00
1	C	426	PRO	CA-CB-CG	-5.11	94.79	104.50

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	7039	0	6874	78	0
1	B	5883	0	5744	93	0
1	C	7213	0	7048	81	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	255	0	212	4	0
2	H	238	0	200	8	0
3	A	126	0	117	3	0
3	B	84	0	78	0	0
3	C	98	0	91	1	0
All	All	20936	0	20364	244	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 6.

All (244) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:CYS:HA	1:A:525:CYS:HB3	1.38	1.05
1:C:65:PHE:O	1:C:264:ALA:HA	1.77	0.85
2:H:97:THR:HG21	2:H:100(B):GLY:HA3	1.59	0.84
2:H:99:ASP:HB2	2:H:100(A):PRO:HD2	1.66	0.77
1:A:358:ILE:HG22	1:A:524:VAL:HG21	1.70	0.74
1:C:815:ARG:NH1	1:C:820:ASP:OD1	2.24	0.71
1:C:902:MET:HB3	1:C:916:LEU:HD11	1.73	0.70
1:C:433:VAL:HG22	1:C:512:VAL:HG22	1.75	0.69
1:A:119:ILE:HG12	1:A:128:ILE:HG13	1.76	0.68
1:A:353:TRP:HZ3	1:A:355:ARG:HB2	1.60	0.67
1:B:231:ILE:HG12	1:B:233:ILE:HG12	1.77	0.66
1:C:903:ALA:HB1	1:C:913:GLN:HG2	1.78	0.66
1:C:395:VAL:HA	1:C:514:SER:O	1.97	0.65
1:C:95:ILE:HG22	1:C:187:LYS:HE2	1.78	0.65
1:A:826:VAL:HB	1:A:1057:PRO:HG2	1.79	0.63
1:A:433:VAL:HG22	1:A:512:VAL:HG22	1.80	0.63
1:B:30:ASN:HB3	1:B:32:PHE:CE1	2.35	0.62
1:B:229:LEU:HD22	1:B:230:PRO:HD2	1.82	0.62
1:B:30:ASN:HB3	1:B:32:PHE:HE1	1.65	0.61
1:B:555:SER:HB2	1:B:586:ASP:HB2	1.82	0.61
1:C:825:LYS:HE2	1:C:939:SER:HA	1.82	0.61
1:A:39:PRO:HG3	1:A:51:THR:HG21	1.83	0.61
1:C:804:GLN:NE2	1:C:935:GLN:OE1	2.23	0.60
1:C:81:ASN:ND2	1:C:240:THR:O	2.33	0.60
1:A:342:PHE:HB2	3:A:1308:NAG:H82	1.83	0.60
1:C:457:ARG:NH1	1:C:459:SER:O	2.34	0.59
1:B:971:GLY:O	1:B:995:ARG:NH1	2.35	0.59
1:C:599:THR:HG22	1:C:601:GLY:H	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ASN:OD1	1:A:123:ALA:N	2.36	0.59
1:A:520:ALA:HB1	1:A:521:PRO:HD3	1.85	0.59
1:B:563:GLN:O	1:B:577:ARG:NH1	2.36	0.58
1:C:318:PHE:HZ	1:C:615:VAL:HG21	1.67	0.58
1:A:1050:MET:HE2	1:A:1052:PHE:CE1	2.39	0.58
1:C:719:THR:HG23	1:C:1068:VAL:HB	1.86	0.58
1:A:811:LYS:NZ	1:A:820:ASP:OD2	2.30	0.58
1:B:64:TRP:HD1	1:B:65:PHE:H	1.50	0.58
1:B:190:ARG:NH2	1:B:264:ALA:O	2.35	0.57
1:B:84:LEU:HD13	1:B:267:VAL:HG21	1.86	0.57
1:B:322:PRO:O	1:B:323:THR:HG22	2.03	0.57
1:A:125:ASN:ND2	1:A:172:SER:O	2.38	0.57
1:A:616:ASN:OD1	1:A:616:ASN:N	2.28	0.57
1:B:104:TRP:HB3	1:B:238:PHE:HE1	1.69	0.57
1:C:46:SER:HA	1:C:279:TYR:O	2.04	0.56
1:A:983:ARG:HD2	1:C:517:LEU:HD13	1.86	0.56
1:B:327:VAL:HG12	1:B:329:PHE:H	1.71	0.55
1:C:418:ILE:HA	1:C:422:ASN:HD22	1.70	0.55
1:C:435:ALA:HB2	1:C:510:VAL:HG13	1.88	0.55
1:C:902:MET:HB3	1:C:916:LEU:CD1	2.36	0.55
2:D:97:THR:HG22	2:D:100(E):LEU:HD23	1.87	0.55
1:B:193:VAL:HB	1:B:204:TYR:HB2	1.88	0.55
1:A:616:ASN:HB3	3:A:1304:NAG:HN2	1.72	0.54
1:A:40:ASP:OD2	1:A:44:ARG:NH2	2.41	0.54
1:B:81:ASN:HD22	1:B:81:ASN:N	2.04	0.54
1:B:298:GLU:OE1	1:B:633:TRP:NE1	2.39	0.54
1:B:957:GLN:NE2	1:C:765:ARG:HD3	2.23	0.54
1:A:562:PHE:HD2	1:B:41:LYS:HD3	1.73	0.54
1:A:662:CYS:HB2	1:A:697:MET:HE3	1.89	0.54
1:B:126:VAL:HG22	1:B:172:SER:HB3	1.89	0.54
1:B:726:ILE:HG13	1:B:1061:VAL:HG22	1.90	0.53
1:C:203:ILE:HB	1:C:227:VAL:HG22	1.89	0.53
1:C:826:VAL:HB	1:C:1057:PRO:HG2	1.90	0.53
1:A:616:ASN:ND2	1:B:836:GLN:HG2	2.23	0.53
1:B:620:VAL:HG13	1:B:621:PRO:HD3	1.91	0.53
1:A:106:PHE:HB2	1:A:117:LEU:HB3	1.90	0.53
1:A:391:CYS:HA	1:A:525:CYS:CB	2.25	0.53
1:B:715:PRO:HA	1:B:1072:GLU:HA	1.90	0.53
1:B:995:ARG:HH22	1:C:991:VAL:HG22	1.73	0.53
1:C:57:PRO:HG3	1:C:273:ARG:HD2	1.89	0.53
1:C:357:ARG:HD3	1:C:396:TYR:CE1	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:980:ILE:HG23	1:A:984:LEU:HD12	1.91	0.52
1:C:629:LEU:HB2	1:C:634:ARG:NH1	2.25	0.52
1:A:659:SER:HB3	1:A:698:SER:HB3	1.91	0.52
1:C:401:VAL:HG22	1:C:509:ARG:HG2	1.91	0.52
1:A:354:ASN:O	1:A:398:ASP:HA	2.10	0.52
1:A:616:ASN:HD21	1:B:836:GLN:HG2	1.75	0.52
1:C:811:LYS:HD2	1:C:812:PRO:HD2	1.92	0.52
1:A:581:THR:HG23	1:A:583:GLU:HG2	1.91	0.51
1:B:1123:SER:OG	1:C:918:GLU:OE2	2.28	0.51
1:C:981:PHE:CZ	1:C:993:ILE:HG13	2.45	0.51
1:B:1127:ASP:OD1	1:B:1127:ASP:N	2.37	0.51
1:C:578:ASP:OD2	1:C:581:THR:OG1	2.23	0.51
1:A:646:ARG:HG3	1:A:646:ARG:HH11	1.75	0.51
1:C:34:ARG:HH21	1:C:191:GLU:CD	2.19	0.51
2:H:30:ASP:O	2:H:52(A):TRP:HB2	2.11	0.51
1:B:202:LYS:HD2	1:B:202:LYS:H	1.76	0.50
1:B:32:PHE:CD2	1:B:218:GLN:HG2	2.46	0.50
1:C:1100:THR:OG1	1:C:1101:HIS:ND1	2.45	0.50
1:B:557:LYS:HD2	1:B:559:PHE:HE1	1.77	0.50
1:B:64:TRP:HD1	1:B:65:PHE:N	2.10	0.49
2:D:99:ASP:HB3	2:D:100(C):TYR:CE2	2.48	0.49
2:H:33:ALA:HB2	2:H:99:ASP:HA	1.94	0.49
2:H:32:TYR:O	2:H:71:ARG:NH2	2.45	0.49
1:B:1107:ARG:HD3	1:C:904:TYR:CZ	2.47	0.49
1:B:91:TYR:CG	1:B:91:TYR:O	2.66	0.49
1:A:497:PHE:CD2	1:A:507:PRO:HB3	2.48	0.48
1:B:194:PHE:HD1	1:B:203:ILE:HG12	1.78	0.48
1:C:726:ILE:HG12	1:C:1061:VAL:HG22	1.96	0.48
1:C:760:CYS:HA	1:C:763:LEU:HD12	1.95	0.48
1:C:354:ASN:O	1:C:398:ASP:HA	2.14	0.48
1:C:364:ASP:OD1	1:C:366:SER:OG	2.22	0.48
1:A:334:ASN:OD1	1:A:334:ASN:N	2.47	0.48
1:A:776:LYS:O	1:A:780:GLU:HG3	2.14	0.48
1:B:960:ASN:HB3	1:B:964:LYS:NZ	2.28	0.48
1:C:715:PRO:HA	1:C:1072:GLU:HA	1.96	0.48
1:A:127:VAL:HG22	1:A:171:VAL:HG22	1.95	0.48
1:C:342:PHE:CE2	1:C:511:VAL:HG11	2.49	0.47
1:C:600:PRO:HB3	1:C:674:TYR:HB2	1.95	0.47
1:A:1079:PRO:HB3	1:B:917:TYR:CE1	2.49	0.47
1:B:65:PHE:CZ	1:B:84:LEU:HD21	2.50	0.47
1:B:90:VAL:HG23	1:B:92:PHE:H	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:LEU:HD21	1:B:119:ILE:HD11	1.94	0.47
1:B:724:THR:HG23	1:B:934:ILE:HD12	1.95	0.47
1:C:426:PRO:HG3	1:C:463:PRO:HB3	1.96	0.47
1:C:1086:LYS:HD2	1:C:1122:VAL:HG11	1.96	0.47
1:B:965:GLN:HE22	1:C:759:PHE:HE2	1.62	0.47
1:A:646:ARG:HG3	1:A:646:ARG:NH1	2.30	0.47
1:A:902:MET:HB3	1:A:916:LEU:HD11	1.96	0.47
1:B:295:PRO:HB2	1:B:608:VAL:HG21	1.96	0.47
1:B:328:ARG:NH1	1:B:531:THR:O	2.47	0.47
1:C:438:SER:HB2	1:C:441:LEU:HB2	1.96	0.47
2:H:97:THR:HG22	2:H:98:TYR:CD1	2.49	0.47
1:A:122:ASN:ND2	1:A:124:THR:OG1	2.47	0.47
1:B:902:MET:HB3	1:B:916:LEU:HD11	1.95	0.47
1:B:960:ASN:HB3	1:B:964:LYS:HZ3	1.79	0.47
1:C:358:ILE:HB	1:C:395:VAL:HG13	1.97	0.46
1:A:298:GLU:HG2	1:A:315:THR:HB	1.97	0.46
1:C:731:MET:HE1	1:C:1014:ARG:HB3	1.96	0.46
1:B:617:CYS:C	1:B:619:GLU:H	2.24	0.46
1:C:616:ASN:HB3	3:C:1304:NAG:HN2	1.80	0.46
1:A:791:THR:HB	1:A:795:LYS:HE2	1.98	0.46
1:B:1045:LYS:HB3	1:B:1045:LYS:HE3	1.65	0.46
2:H:100:SER:N	2:H:100(A):PRO:HD2	2.30	0.46
1:C:172:SER:OG	1:C:173:GLN:N	2.49	0.46
1:B:310:LYS:HG3	1:B:600:PRO:HA	1.98	0.45
1:C:630:THR:O	1:C:632:THR:HG23	2.16	0.45
1:A:758:SER:O	1:A:762:GLN:HG3	2.16	0.45
1:B:227:VAL:HG22	1:B:228:ASP:H	1.82	0.45
1:A:206:LYS:HB2	1:A:223:LEU:HA	1.98	0.45
1:B:555:SER:HB2	1:B:586:ASP:CB	2.46	0.45
1:B:953:ASN:O	1:B:957:GLN:HB3	2.17	0.45
1:A:344:ALA:HB3	1:A:347:PHE:CE1	2.52	0.45
1:B:555:SER:OG	1:B:557:LYS:HG3	2.17	0.45
1:A:1127:ASP:OD1	1:A:1127:ASP:N	2.49	0.45
1:B:551:VAL:HG12	1:B:553:THR:HG23	1.98	0.45
1:C:785:VAL:HG21	1:C:789:TYR:HE2	1.81	0.45
1:C:1090:PRO:HD3	1:C:1095:PHE:CE2	2.52	0.45
1:B:620:VAL:N	1:B:621:PRO:HD2	2.32	0.45
1:B:825:LYS:HD2	1:B:825:LYS:HA	1.88	0.45
1:A:34:ARG:HH21	1:A:191:GLU:CD	2.23	0.44
1:A:1106:GLN:HG3	1:A:1109:PHE:O	2.17	0.44
1:A:105:ILE:HG22	1:A:118:LEU:HD23	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:599:THR:HG22	1:A:608:VAL:HG12	1.99	0.44
1:C:295:PRO:HG3	1:C:633:TRP:CZ3	2.53	0.44
1:A:917:TYR:CE1	1:C:1079:PRO:HB3	2.52	0.44
1:A:193:VAL:HG13	1:A:270:LEU:HD11	1.99	0.44
1:B:190:ARG:O	1:B:190:ARG:HG2	2.17	0.44
1:A:42:VAL:HG21	1:C:567:ARG:HH21	1.81	0.44
1:B:552:LEU:HA	1:B:586:ASP:O	2.18	0.44
1:B:592:PHE:CD1	1:C:740:MET:HE1	2.53	0.44
1:C:119:ILE:HG12	1:C:128:ILE:HG12	2.00	0.44
1:B:119:ILE:HG22	1:B:121:ASN:OD1	2.18	0.44
1:B:64:TRP:CD1	1:B:65:PHE:N	2.86	0.43
1:B:802:PHE:HZ	1:B:898:PHE:CZ	2.36	0.43
1:C:103:GLY:O	1:C:241:LEU:N	2.44	0.43
1:A:521:PRO:HD2	1:A:521:PRO:O	2.18	0.43
1:C:659:SER:HB3	1:C:698:SER:HB3	2.00	0.43
1:A:127:VAL:C	1:A:128:ILE:HD12	2.43	0.43
2:D:99:ASP:HB3	2:D:100(C):TYR:HE2	1.82	0.43
1:B:328:ARG:HH21	1:B:578:ASP:CG	2.26	0.43
1:B:708:SER:HB3	1:B:711:SER:HB3	2.00	0.43
1:B:32:PHE:N	1:B:32:PHE:CD1	2.86	0.43
1:B:328:ARG:HG3	1:B:543:PHE:CE2	2.53	0.43
1:B:555:SER:H	1:B:586:ASP:H	1.66	0.43
1:A:623:ALA:HB3	1:A:634:ARG:HH12	1.83	0.43
1:B:81:ASN:N	1:B:81:ASN:ND2	2.66	0.43
1:B:117:LEU:HD13	1:B:231:ILE:HD12	2.01	0.43
1:A:172:SER:OG	1:A:173:GLN:N	2.49	0.43
1:A:281:GLU:N	1:A:281:GLU:OE1	2.51	0.43
1:A:36:VAL:O	1:A:222:ALA:HA	2.19	0.43
3:A:1304:NAG:H82	1:B:834:ILE:HG21	2.00	0.43
1:A:631:PRO:O	1:A:632:THR:HG22	2.19	0.42
1:A:804:GLN:OE1	1:A:935:GLN:NE2	2.39	0.42
1:A:1088:HIS:HB3	1:A:1120:THR:HG21	2.01	0.42
1:B:715:PRO:HD3	1:C:894:LEU:HD13	2.01	0.42
1:A:722:VAL:HG22	1:A:1065:VAL:HG22	2.01	0.42
1:C:39:PRO:HG3	1:C:51:THR:HG21	2.00	0.42
1:C:426:PRO:HD2	1:C:429:PHE:HB2	1.99	0.42
1:A:1088:HIS:HB3	1:A:1120:THR:CG2	2.48	0.42
1:A:983:ARG:O	1:C:382:VAL:HA	2.19	0.42
1:B:129:LYS:HG2	1:B:169:GLU:HG2	2.01	0.42
1:C:612:TYR:HE2	1:C:651:ILE:HD12	1.85	0.42
1:A:240:THR:HG21	1:A:265:TYR:CZ	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:LYS:HD3	1:B:204:TYR:HE1	1.85	0.42
1:B:959:LEU:HD23	1:B:959:LEU:HA	1.92	0.42
1:C:40:ASP:OD2	1:C:44:ARG:NH2	2.51	0.42
1:A:900:MET:HE3	1:A:900:MET:HB2	1.77	0.42
2:D:100(A):PRO:HA	2:D:100(F):ASN:ND2	2.35	0.42
1:B:220:PHE:HE2	1:B:285:ILE:HG22	1.84	0.42
1:B:722:VAL:HG22	1:B:1065:VAL:HG22	2.02	0.42
1:A:64:TRP:CD1	1:A:64:TRP:C	2.98	0.42
1:B:900:MET:HE3	1:B:900:MET:HB3	1.76	0.42
1:A:318:PHE:HA	1:A:631:PRO:HD3	2.02	0.41
1:A:894:LEU:HD13	1:C:715:PRO:HD3	2.01	0.41
1:B:996:LEU:HD23	1:B:996:LEU:HA	1.92	0.41
1:B:1106:GLN:HG3	1:B:1109:PHE:O	2.19	0.41
1:A:623:ALA:HB3	1:A:634:ARG:NH1	2.35	0.41
1:C:577:ARG:HG3	1:C:577:ARG:HH11	1.83	0.41
1:C:900:MET:HE3	1:C:900:MET:HB3	1.79	0.41
1:C:979:ASP:O	1:C:983:ARG:HB2	2.21	0.41
1:A:600:PRO:HB3	1:A:674:TYR:HB2	2.03	0.41
1:B:578:ASP:OD2	1:B:581:THR:HG22	2.21	0.41
1:B:745:ASP:OD1	1:B:745:ASP:C	2.63	0.41
1:B:1079:PRO:HB3	1:C:917:TYR:CE1	2.56	0.41
1:C:616:ASN:OD1	1:C:616:ASN:N	2.52	0.41
1:A:41:LYS:HD2	1:C:562:PHE:HD1	1.85	0.41
1:A:962:LEU:HD12	1:A:962:LEU:HA	1.94	0.41
1:C:129:LYS:HG2	1:C:169:GLU:HG2	2.03	0.41
1:A:902:MET:HB3	1:A:916:LEU:CD1	2.51	0.41
1:B:804:GLN:OE1	1:B:935:GLN:HG2	2.21	0.41
1:C:403:ARG:HB2	1:C:495:TYR:CD1	2.56	0.41
1:B:106:PHE:HB3	1:B:235:ILE:HG23	2.03	0.41
1:B:1079:PRO:HG2	1:B:1130:ILE:O	2.21	0.41
1:B:1125:ASN:HB2	1:B:1127:ASP:OD1	2.20	0.41
1:C:442:ASP:OD1	2:H:98:TYR:OH	2.29	0.41
1:C:442:ASP:OD1	1:C:451:TYR:OH	2.27	0.41
1:B:763:LEU:HD22	1:B:1008:VAL:HG21	2.03	0.41
1:A:353:TRP:CZ3	1:A:355:ARG:HB2	2.47	0.40
1:A:377:PHE:C	1:A:377:PHE:CD2	2.98	0.40
1:A:849:LEU:HD12	1:A:849:LEU:HA	1.96	0.40
1:B:703:ASN:OD1	1:B:703:ASN:C	2.64	0.40
1:C:416:GLY:O	1:C:420:ASP:HB2	2.21	0.40
1:C:1082:CYS:HB2	1:C:1132:ILE:HG12	2.03	0.40
1:A:226:LEU:HD23	1:A:226:LEU:HA	1.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:965:GLN:HE22	1:B:762:GLN:HE22	1.68	0.40
1:B:27:ALA:O	1:B:64:TRP:N	2.52	0.40
1:C:802:PHE:HZ	1:C:898:PHE:CZ	2.39	0.40
1:B:633:TRP:HE3	1:B:633:TRP:O	2.03	0.40
1:C:298:GLU:HB3	1:C:315:THR:HG21	2.03	0.40
1:A:118:LEU:HB3	1:A:129:LYS:HG3	2.03	0.40
1:A:357:ARG:C	1:A:358:ILE:HD12	2.46	0.40
1:B:32:PHE:N	1:B:32:PHE:HD1	2.18	0.40
1:B:47:VAL:HG12	1:B:279:TYR:HB2	2.04	0.40
1:B:878:LEU:HD21	1:B:1052:PHE:HB3	2.04	0.40
1:C:128:ILE:HD13	1:C:229:LEU:HD11	2.04	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	888/1249 (71%)	862 (97%)	24 (3%)	2 (0%)	43	61
1	B	735/1249 (59%)	703 (96%)	29 (4%)	3 (0%)	30	46
1	C	902/1249 (72%)	866 (96%)	34 (4%)	2 (0%)	43	61
2	D	22/449 (5%)	21 (96%)	1 (4%)	0	100	100
2	H	21/449 (5%)	17 (81%)	4 (19%)	0	100	100
All	All	2568/4645 (55%)	2469 (96%)	92 (4%)	7 (0%)	37	53

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	557	LYS
1	C	231	ILE
1	C	446	SER

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Mol	Chain	Res	Type
1	A	519	HIS
1	B	592	PHE
1	B	189	LEU
1	A	520	ALA

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	776/1094 (71%)	757 (98%)	19 (2%)	43	67
1	B	645/1094 (59%)	634 (98%)	11 (2%)	53	75
1	C	792/1094 (72%)	778 (98%)	14 (2%)	51	74
2	D	28/387 (7%)	28 (100%)	0	100	100
2	H	25/387 (6%)	25 (100%)	0	100	100
All	All	2266/4056 (56%)	2222 (98%)	44 (2%)	49	73

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	LEU
1	A	118	LEU
1	A	127	VAL
1	A	281	GLU
1	A	349	SER
1	A	383	SER
1	A	394	ASN
1	A	430	THR
1	A	434	ILE
1	A	555	SER
1	A	588	THR
1	A	616	ASN
1	A	620	VAL
1	A	630	THR
1	A	632	THR

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Mol	Chain	Res	Type
1	A	791	THR
1	A	900	MET
1	A	1120	THR
1	A	1133	VAL
1	B	81	ASN
1	B	240	THR
1	B	284	THR
1	B	539	VAL
1	B	620	VAL
1	B	633	TRP
1	B	747	THR
1	B	791	THR
1	B	965	GLN
1	B	1100	THR
1	B	1136	THR
1	C	46	SER
1	C	307	THR
1	C	324	GLU
1	C	399	SER
1	C	407	VAL
1	C	455	LEU
1	C	531	THR
1	C	615	VAL
1	C	630	THR
1	C	709	ASN
1	C	747	THR
1	C	754	LEU
1	C	1125	ASN
1	C	1136	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	122	ASN
1	A	239	GLN
1	A	314	GLN
1	A	505	HIS
1	A	755	GLN
1	A	965	GLN
1	B	196	ASN
1	B	644	GLN

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Mol	Chain	Res	Type
1	B	658	ASN
1	B	1071	GLN
1	B	1125	ASN
1	C	66	HIS
1	C	81	ASN
1	C	450	ASN
1	C	519	HIS
1	C	965	GLN
1	C	1002	GLN
1	C	1005	GLN
1	C	1125	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

22 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	A	1308	1	14,14,15	0.26	0	17,19,21	0.50	0
3	NAG	A	1305	1	14,14,15	0.28	0	17,19,21	0.37	0
3	NAG	A	1301	1	14,14,15	0.22	0	17,19,21	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	A	1307	1	14,14,15	0.54	0	17,19,21	0.71	1 (5%)
3	NAG	C	1304	1	14,14,15	0.24	0	17,19,21	0.45	0
3	NAG	A	1306	1	14,14,15	0.28	0	17,19,21	0.53	0
3	NAG	C	1302	1	14,14,15	0.21	0	17,19,21	0.45	0
3	NAG	C	1306	1	14,14,15	0.30	0	17,19,21	0.51	0
3	NAG	C	1307	1	14,14,15	0.26	0	17,19,21	0.50	0
3	NAG	B	1306	1	14,14,15	0.27	0	17,19,21	0.51	0
3	NAG	C	1301	1	14,14,15	0.23	0	17,19,21	0.54	0
3	NAG	B	1305	1	14,14,15	0.29	0	17,19,21	0.51	0
3	NAG	C	1303	1	14,14,15	0.53	0	17,19,21	0.63	0
3	NAG	C	1305	1	14,14,15	0.32	0	17,19,21	0.44	0
3	NAG	B	1303	1	14,14,15	0.24	0	17,19,21	0.41	0
3	NAG	A	1302	1	14,14,15	0.18	0	17,19,21	0.43	0
3	NAG	A	1304	1	14,14,15	0.36	0	17,19,21	0.51	0
3	NAG	A	1309	1	14,14,15	0.24	0	17,19,21	0.54	0
3	NAG	B	1301	1	14,14,15	0.26	0	17,19,21	0.51	0
3	NAG	B	1302	1	14,14,15	0.19	0	17,19,21	0.44	0
3	NAG	A	1303	1	14,14,15	0.28	0	17,19,21	0.42	0
3	NAG	B	1304	1	14,14,15	0.19	0	17,19,21	0.53	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
3	NAG	A	1308	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1305	1	-	1/6/23/26	0/1/1/1
3	NAG	A	1301	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1307	1	-	4/6/23/26	0/1/1/1
3	NAG	C	1304	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1306	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1302	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1306	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1307	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1306	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1301	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1305	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1303	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1305	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	1303	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1302	1	-	1/6/23/26	0/1/1/1
3	NAG	A	1304	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1309	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1301	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1302	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1303	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1304	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1307	NAG	C1-O5-C5	2.36	115.35	112.19

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1307	NAG	C4-C5-C6-O6
3	C	1302	NAG	O5-C5-C6-O6
3	C	1303	NAG	O5-C5-C6-O6
3	A	1301	NAG	O5-C5-C6-O6
3	A	1303	NAG	O5-C5-C6-O6
3	A	1303	NAG	C4-C5-C6-O6
3	B	1303	NAG	O5-C5-C6-O6
3	C	1303	NAG	C4-C5-C6-O6
3	B	1304	NAG	O5-C5-C6-O6
3	A	1307	NAG	O5-C5-C6-O6
3	C	1305	NAG	O5-C5-C6-O6
3	A	1308	NAG	O5-C5-C6-O6
3	C	1302	NAG	C4-C5-C6-O6
3	A	1301	NAG	C4-C5-C6-O6
3	C	1305	NAG	C4-C5-C6-O6
3	A	1308	NAG	C4-C5-C6-O6
3	B	1303	NAG	C4-C5-C6-O6
3	B	1304	NAG	C4-C5-C6-O6
3	A	1307	NAG	C8-C7-N2-C2
3	A	1307	NAG	O7-C7-N2-C2
3	B	1302	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	C	1301	NAG	O5-C5-C6-O6
3	C	1301	NAG	C4-C5-C6-O6
3	B	1305	NAG	O5-C5-C6-O6
3	B	1305	NAG	C4-C5-C6-O6
3	A	1302	NAG	O5-C5-C6-O6
3	B	1302	NAG	C4-C5-C6-O6
3	A	1306	NAG	C4-C5-C6-O6
3	A	1306	NAG	O5-C5-C6-O6
3	A	1305	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1308	NAG	1	0
3	C	1304	NAG	1	0
3	A	1304	NAG	2	0

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.

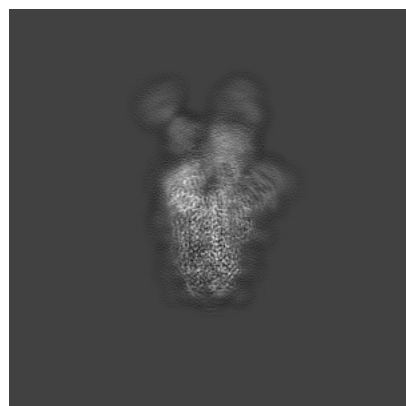
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34469. These allow visual inspection of the internal detail of the map and identification of artifacts.

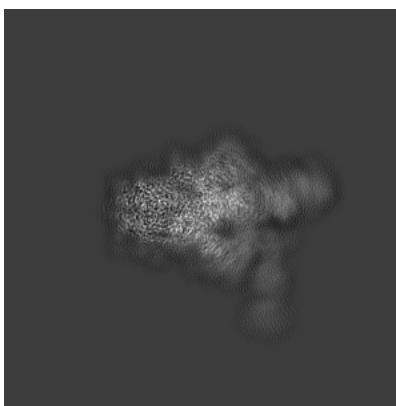
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

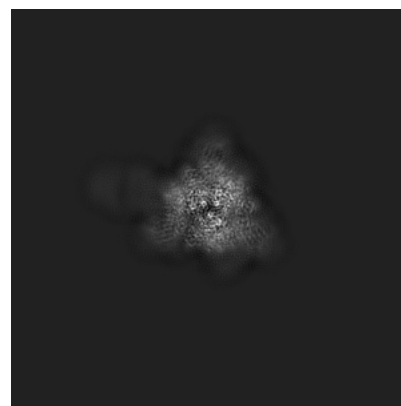
6.1.1 Primary map



X

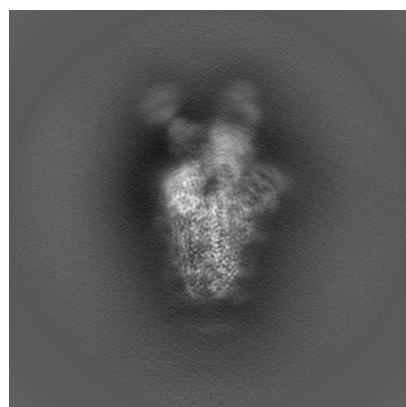


Y

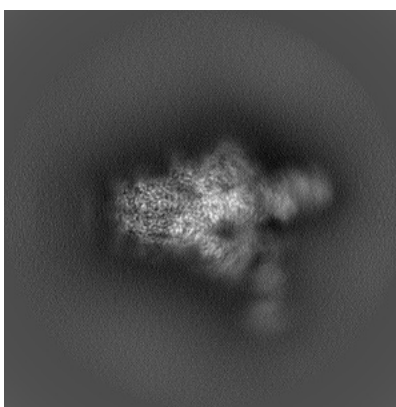


Z

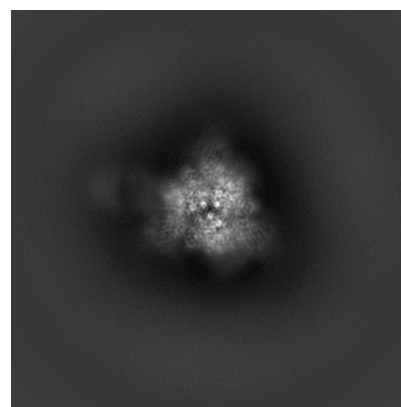
6.1.2 Raw map



X



Y

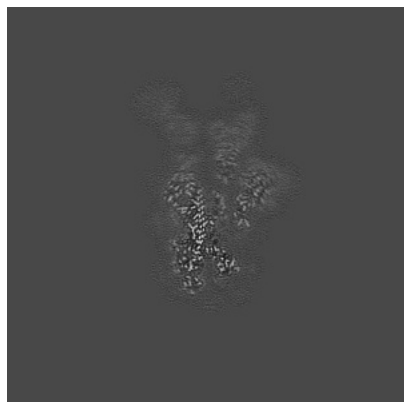


Z

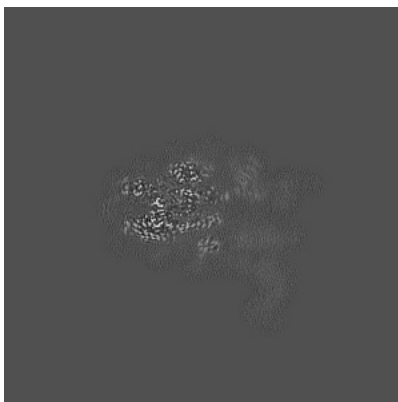
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

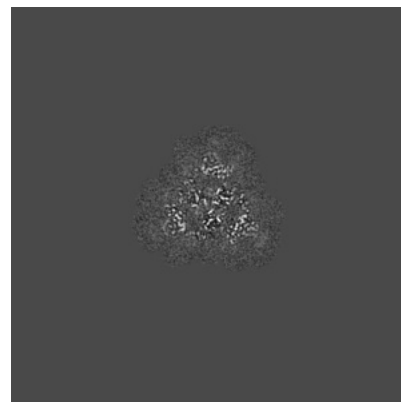
6.2.1 Primary map



X Index: 256

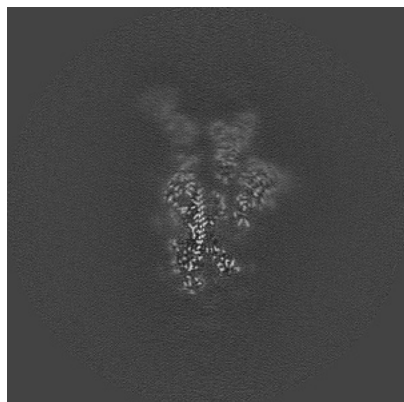


Y Index: 256

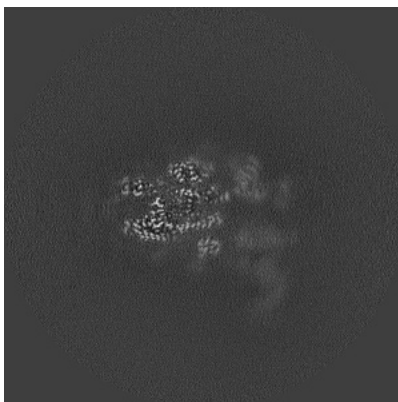


Z Index: 256

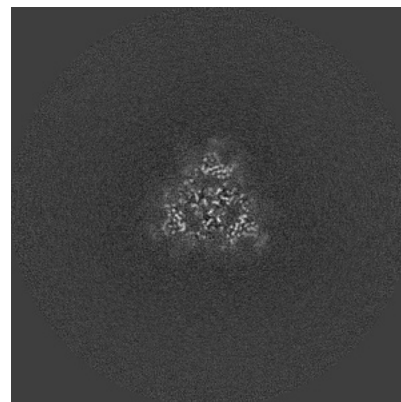
6.2.2 Raw map



X Index: 256



Y Index: 256

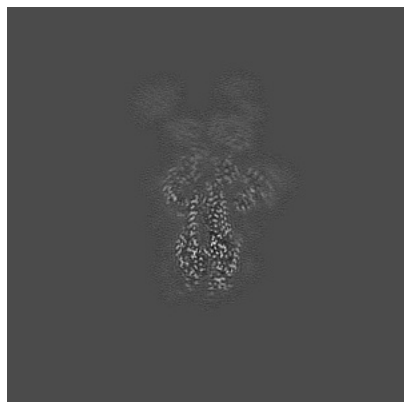


Z Index: 256

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

6.3 Largest variance slices [i](#)

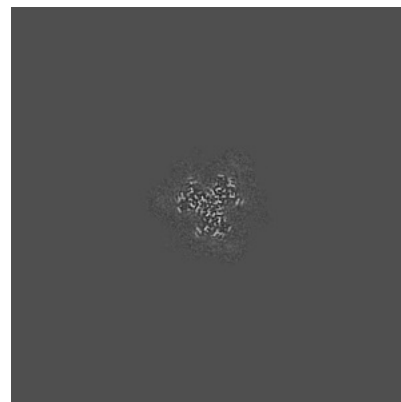
6.3.1 Primary map



X Index: 267

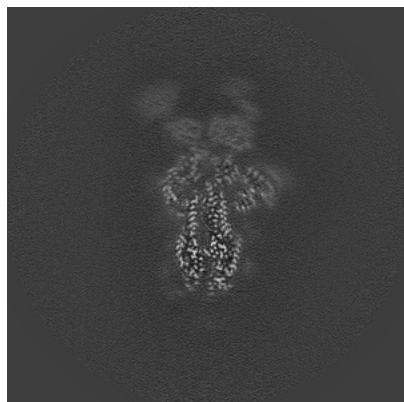


Y Index: 265

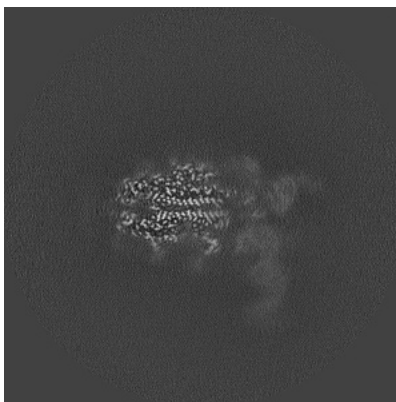


Z Index: 199

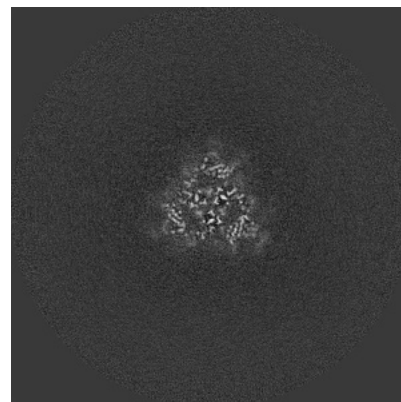
6.3.2 Raw map



X Index: 267



Y Index: 264

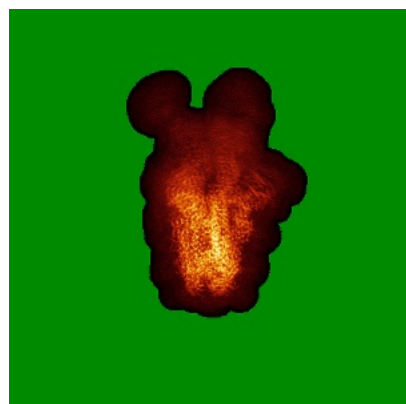


Z Index: 255

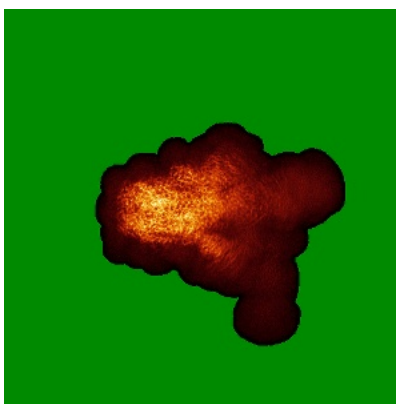
The images above show the largest variance slices of the map in three orthogonal directions.

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

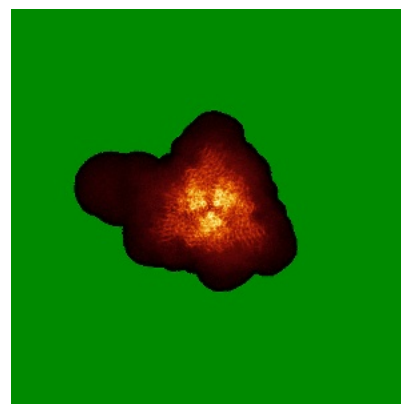
6.4.1 Primary map



X

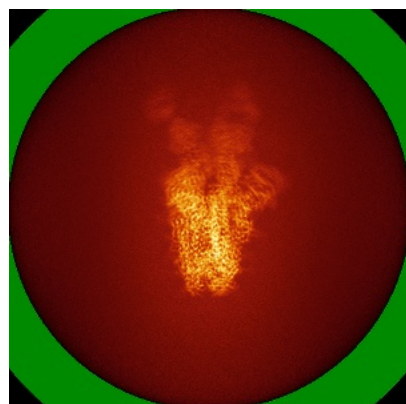


Y

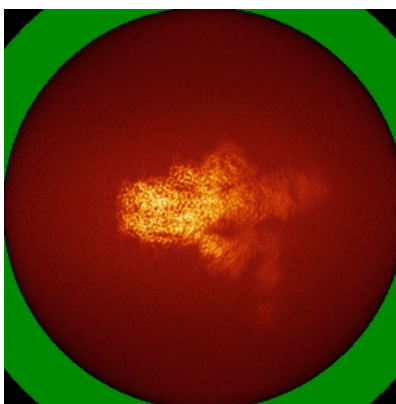


Z

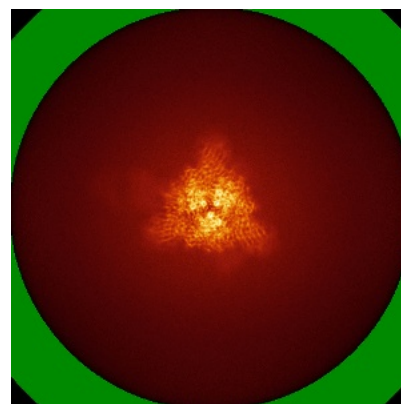
6.4.2 Raw map



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.

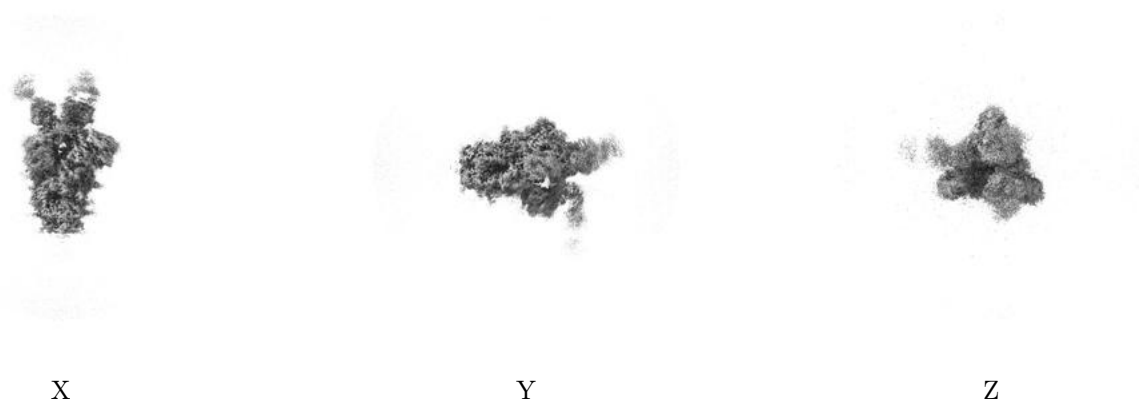
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.005. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

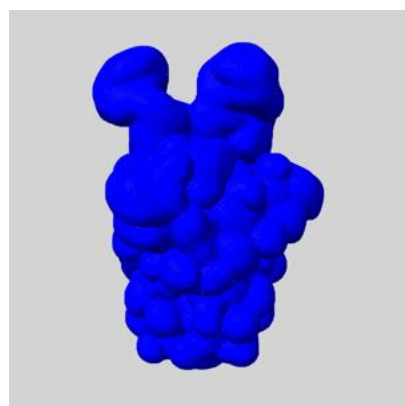
6.6 Mask visualisation [i](#)

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

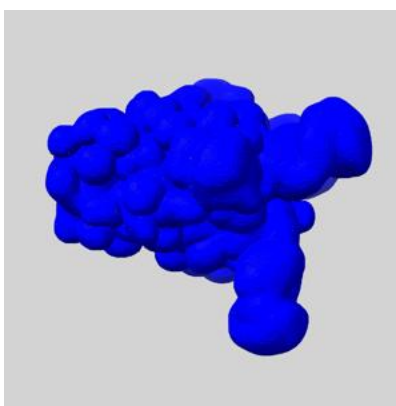
A mask typically either:

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

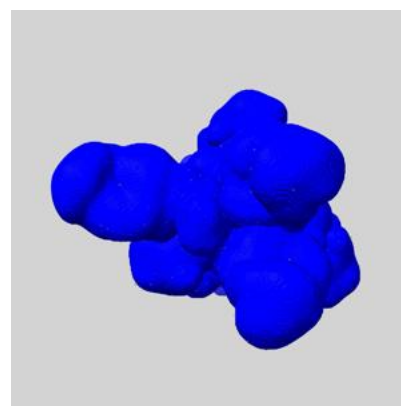
6.6.1 emd_34469_msk_1.map [i](#)



X



Y

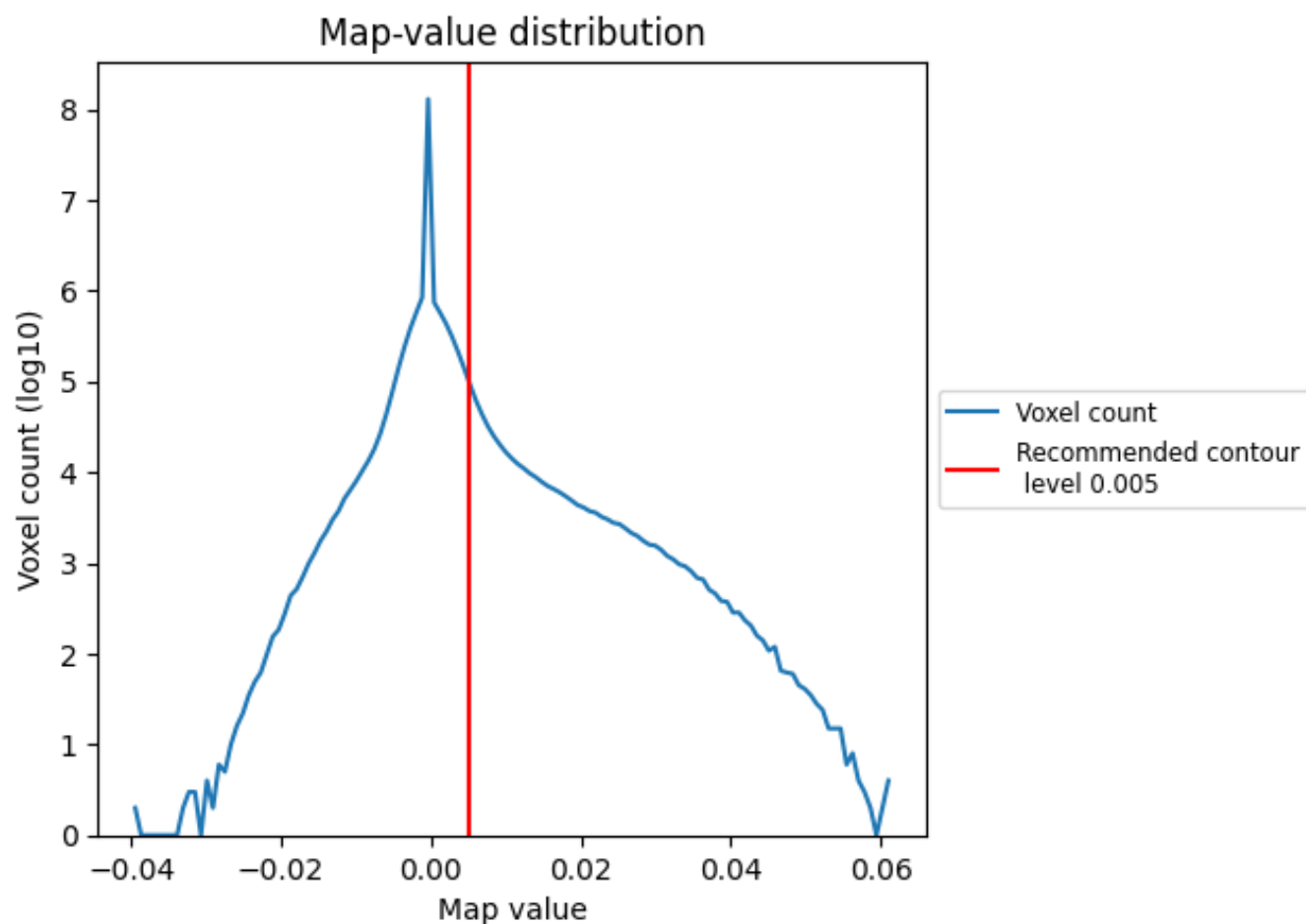


Z

7 Map analysis [i](#)

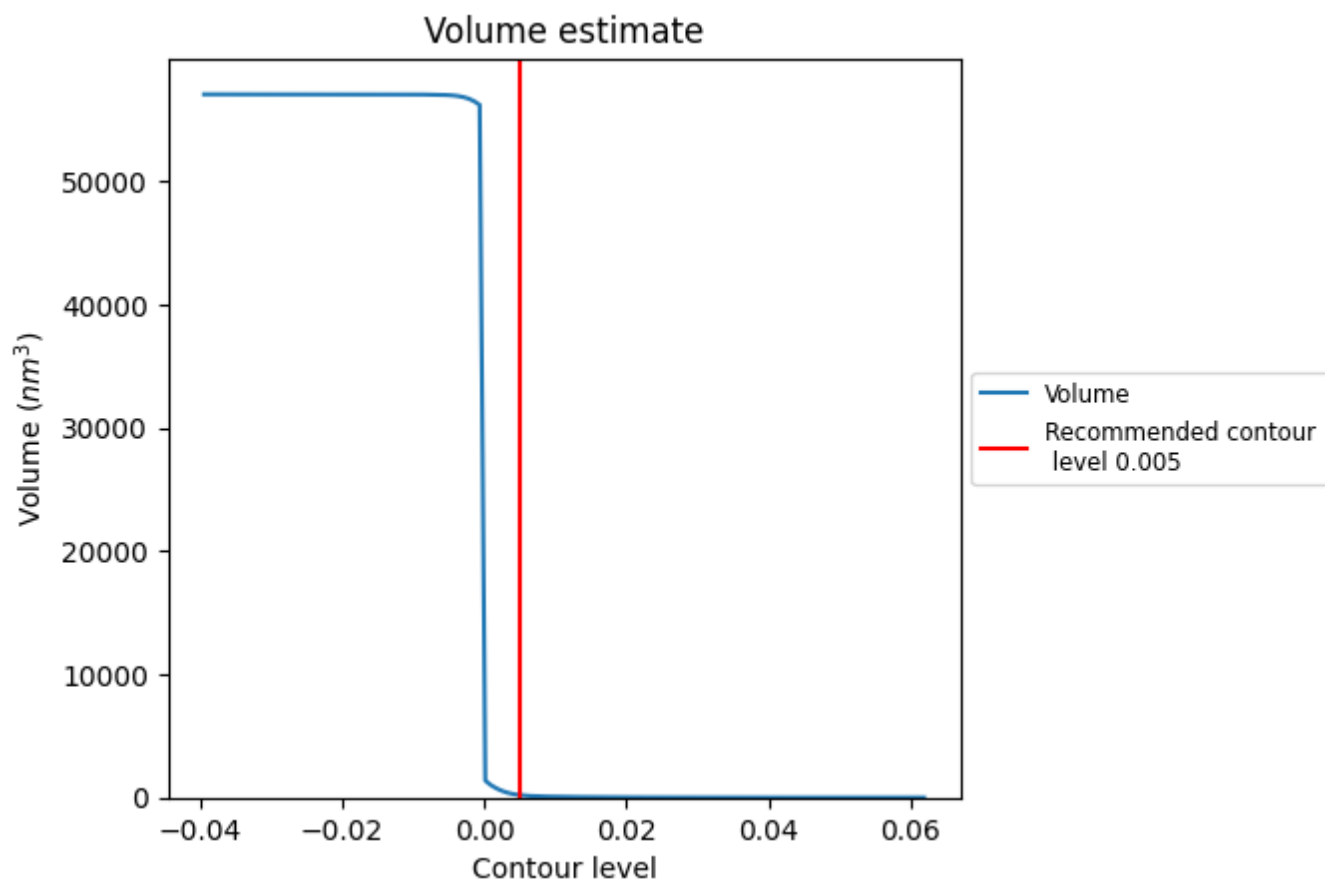
This section contains the results of statistical analysis of the map.

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.

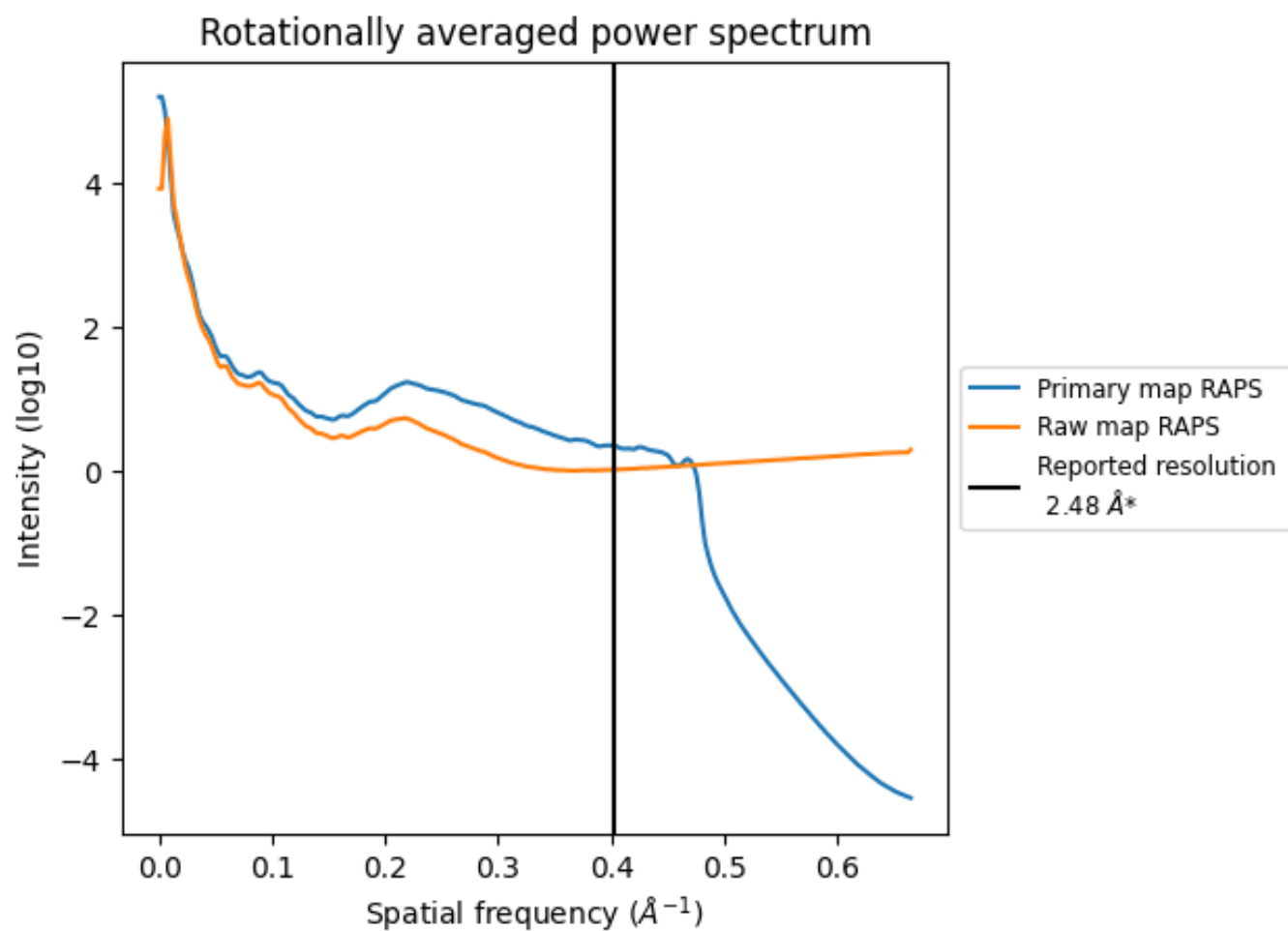
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 201 nm³; this corresponds to an approximate mass of 181 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 ⓘ

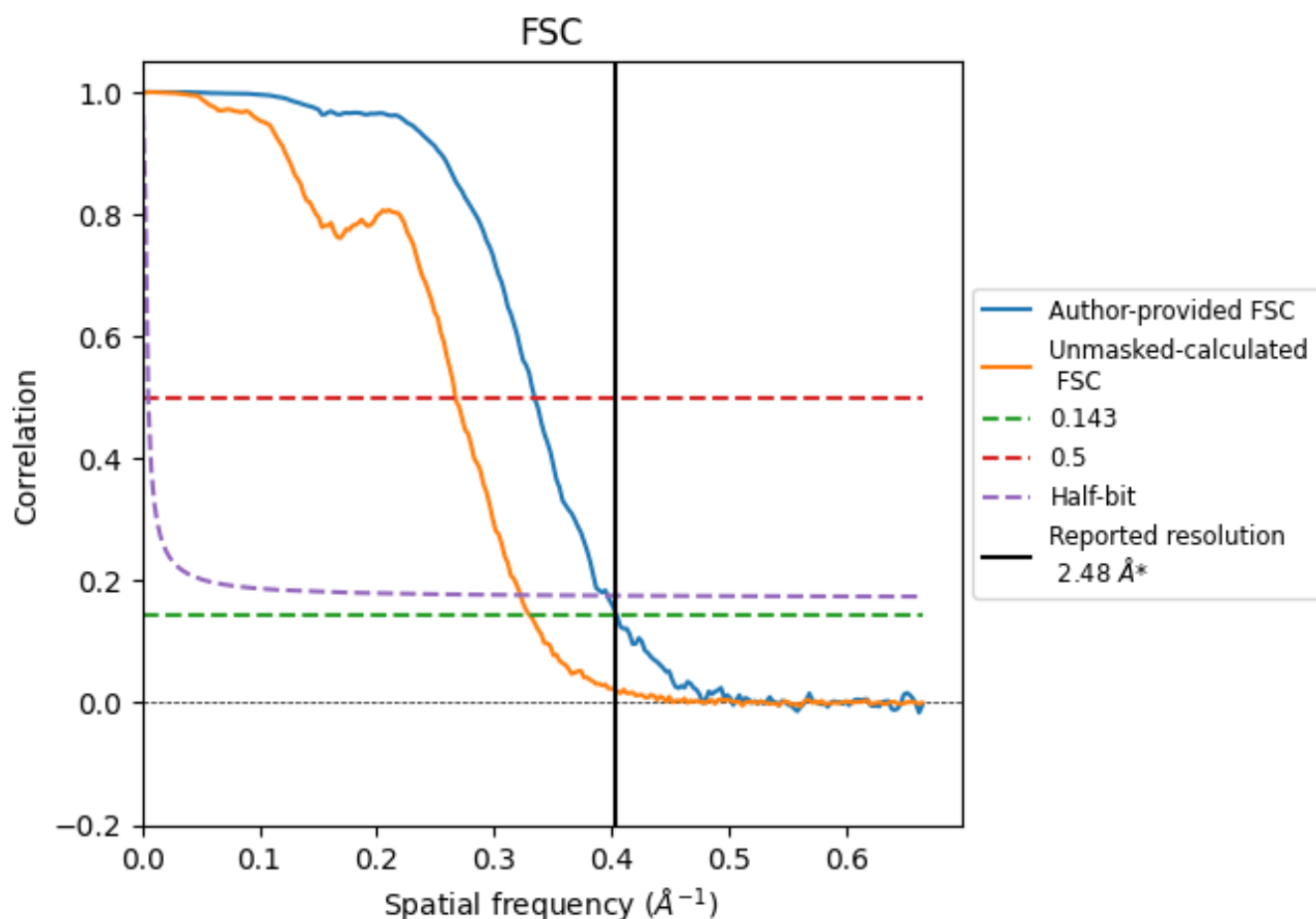


*Reported resolution corresponds to spatial frequency of 0.403 Å⁻¹

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 $0.403 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

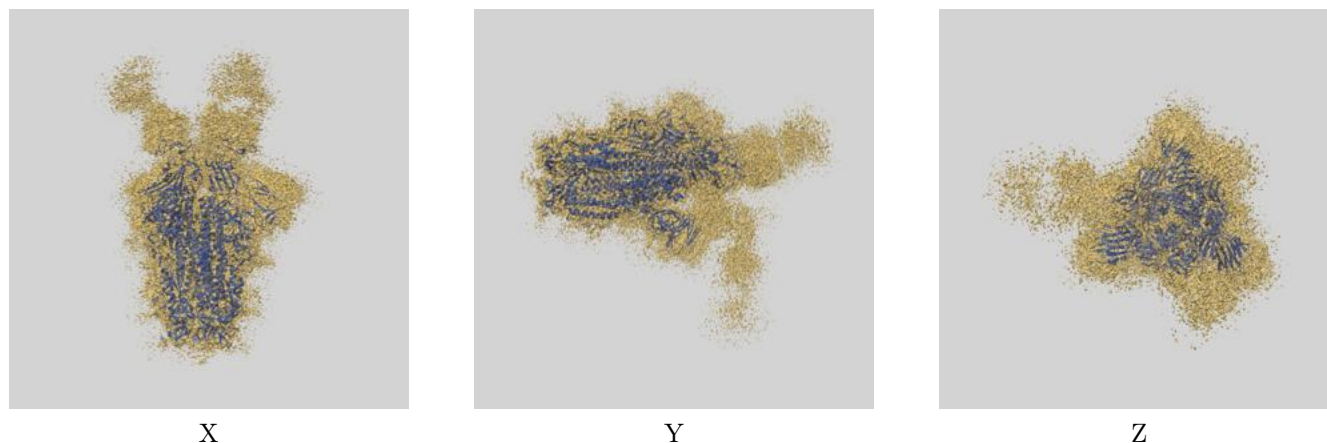
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.48	-	-
Author-provided FSC curve	2.47	2.99	2.52
Unmasked-calculated*	3.03	3.74	3.10

*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 3.03 differs from the reported value 2.48 by more than 10 %

9 Map-model fit [i](#)

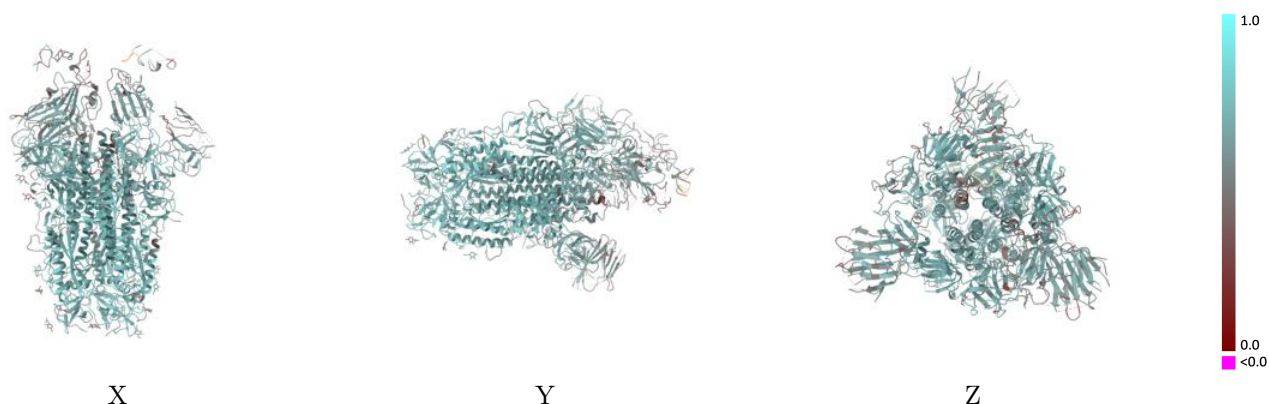
This section contains information regarding the fit between EMDB map EMD-34469 and PDB model 8H3M. Per-residue inclusion information can be found in section 3 on page 13.

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.005 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](#)



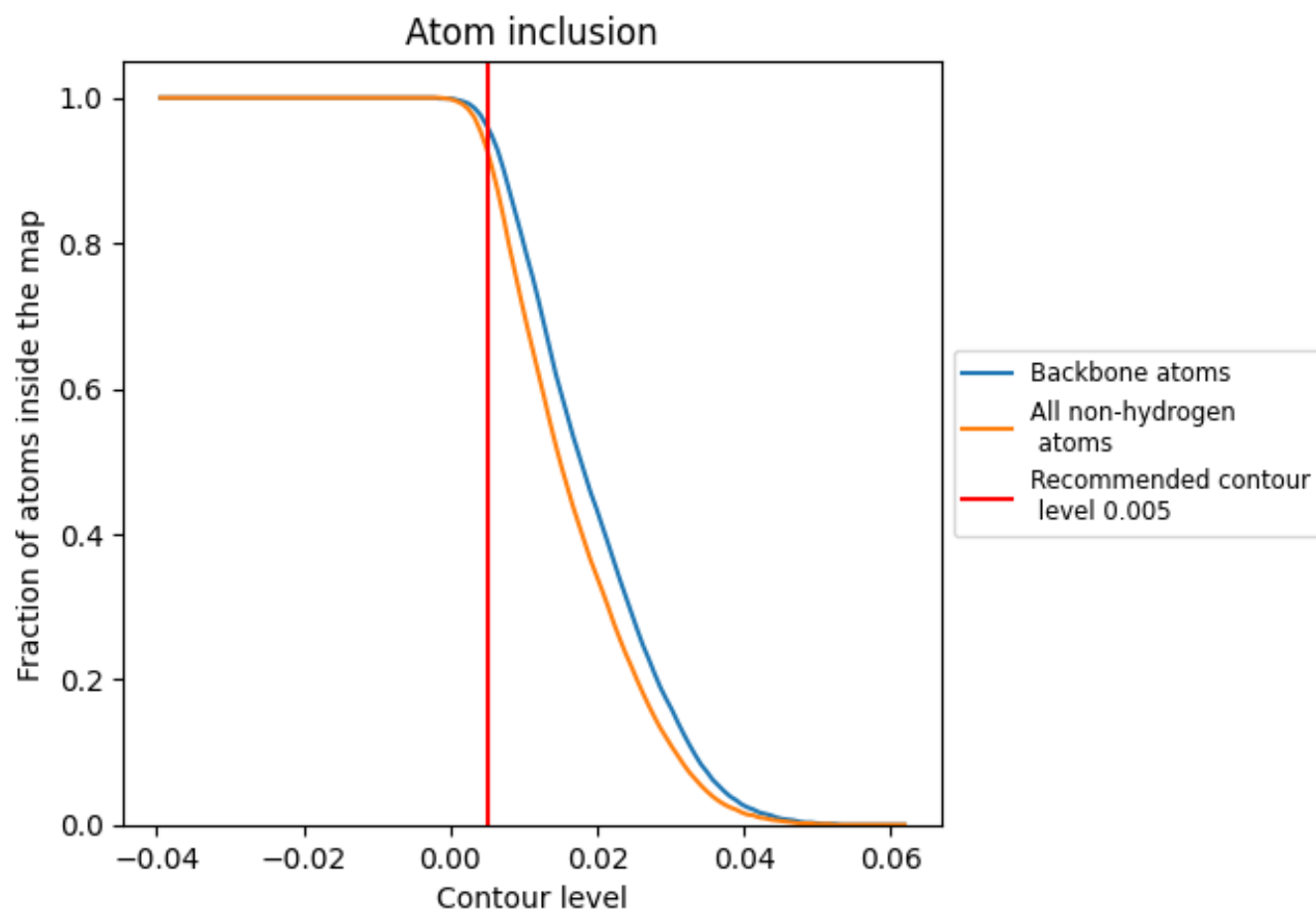
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.005).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% 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 (0.005) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9290	0.6350
A	0.9390	0.6430
B	0.9430	0.6470
C	0.9250	0.6300
D	0.6810	0.4580
H	0.6720	0.4540

