



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 10:27 AM UTC

PDB ID : 7KXJ / pdb_00007kxj
EMDB ID : EMD-23064
Title : SARS-CoV-2 spike protein in complex with Fab 15033-7, 3-"up", asymmetric
Authors : Li, Z.; Rini, J.
Deposited on : 2020-12-04
Resolution : 6.40 Å(reported)
Based on initial model : 7KML

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

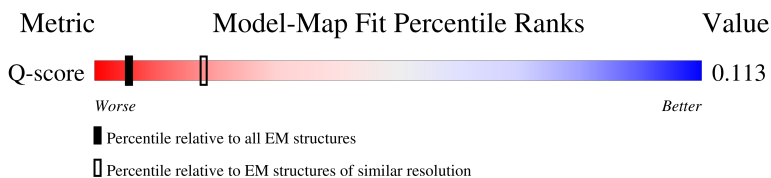
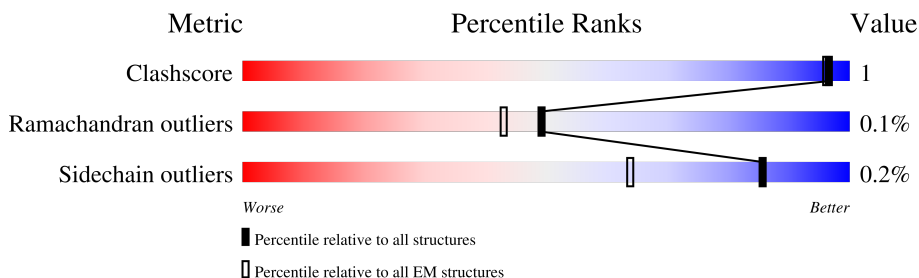
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

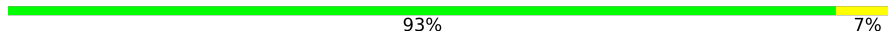
The reported resolution of this entry is 6.40 Å.

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.



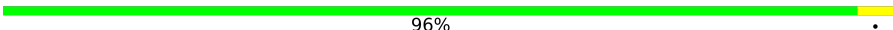
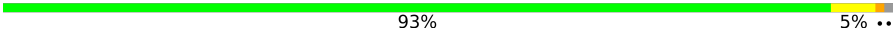
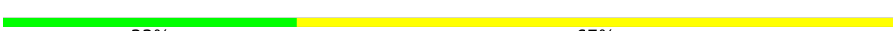
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	544 (5.90 - 6.90)

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	1274	
1	B	1274	
1	C	1274	
2	L	214	

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Mol	Chain	Length	Quality of chain
2	M	214	 96% .
2	N	214	 21% 89% 10%
3	H	225	 93% 5% ..
3	I	225	 92% 7% .
3	J	225	 39% 92% 6% .
4	D	3	 67% 33%
4	E	3	 33% 67%
5	F	2	 50% 50%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 34083 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	1026	Total	C	N	O	S	0	0
			8017	5118	1336	1525	38		
1	B	1026	Total	C	N	O	S	0	0
			8017	5118	1336	1525	38		
1	C	1026	Total	C	N	O	S	0	0
			8017	5118	1336	1525	38		

There are 216 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	682	SER	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	1212	SER	-	expression tag	UNP P0DTC2
A	1213	GLY	-	expression tag	UNP P0DTC2
A	1214	GLY	-	expression tag	UNP P0DTC2
A	1215	TYR	-	expression tag	UNP P0DTC2
A	1216	ILE	-	expression tag	UNP P0DTC2
A	1217	PRO	-	expression tag	UNP P0DTC2
A	1218	GLU	-	expression tag	UNP P0DTC2
A	1219	ALA	-	expression tag	UNP P0DTC2
A	1220	PRO	-	expression tag	UNP P0DTC2
A	1221	ARG	-	expression tag	UNP P0DTC2
A	1222	ASP	-	expression tag	UNP P0DTC2
A	1223	GLY	-	expression tag	UNP P0DTC2
A	1224	GLN	-	expression tag	UNP P0DTC2
A	1225	ALA	-	expression tag	UNP P0DTC2
A	1226	TYR	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1227	VAL	-	expression tag	UNP P0DTC2
A	1228	ARG	-	expression tag	UNP P0DTC2
A	1229	LYS	-	expression tag	UNP P0DTC2
A	1230	ASP	-	expression tag	UNP P0DTC2
A	1231	GLY	-	expression tag	UNP P0DTC2
A	1232	GLU	-	expression tag	UNP P0DTC2
A	1233	TRP	-	expression tag	UNP P0DTC2
A	1234	VAL	-	expression tag	UNP P0DTC2
A	1235	LEU	-	expression tag	UNP P0DTC2
A	1236	LEU	-	expression tag	UNP P0DTC2
A	1237	SER	-	expression tag	UNP P0DTC2
A	1238	THR	-	expression tag	UNP P0DTC2
A	1239	PHE	-	expression tag	UNP P0DTC2
A	1240	LEU	-	expression tag	UNP P0DTC2
A	1241	ASN	-	expression tag	UNP P0DTC2
A	1242	SER	-	expression tag	UNP P0DTC2
A	1243	GLY	-	expression tag	UNP P0DTC2
A	1244	ARG	-	expression tag	UNP P0DTC2
A	1245	ALA	-	expression tag	UNP P0DTC2
A	1246	HIS	-	expression tag	UNP P0DTC2
A	1247	HIS	-	expression tag	UNP P0DTC2
A	1248	HIS	-	expression tag	UNP P0DTC2
A	1249	HIS	-	expression tag	UNP P0DTC2
A	1250	HIS	-	expression tag	UNP P0DTC2
A	1251	HIS	-	expression tag	UNP P0DTC2
A	1252	GLY	-	expression tag	UNP P0DTC2
A	1253	ALA	-	expression tag	UNP P0DTC2
A	1254	GLY	-	expression tag	UNP P0DTC2
A	1255	GLY	-	expression tag	UNP P0DTC2
A	1256	LEU	-	expression tag	UNP P0DTC2
A	1257	ASN	-	expression tag	UNP P0DTC2
A	1258	ASP	-	expression tag	UNP P0DTC2
A	1259	ILE	-	expression tag	UNP P0DTC2
A	1260	PHE	-	expression tag	UNP P0DTC2
A	1261	GLU	-	expression tag	UNP P0DTC2
A	1262	ALA	-	expression tag	UNP P0DTC2
A	1263	GLN	-	expression tag	UNP P0DTC2
A	1264	LYS	-	expression tag	UNP P0DTC2
A	1265	ILE	-	expression tag	UNP P0DTC2
A	1266	GLU	-	expression tag	UNP P0DTC2
A	1267	TRP	-	expression tag	UNP P0DTC2
A	1268	HIS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1269	GLU	-	expression tag	UNP P0DTC2
A	1270	ASP	-	expression tag	UNP P0DTC2
A	1271	THR	-	expression tag	UNP P0DTC2
A	1272	ALA	-	expression tag	UNP P0DTC2
A	1273	ALA	-	expression tag	UNP P0DTC2
A	1274	ALA	-	expression tag	UNP P0DTC2
B	682	SER	ARG	conflict	UNP P0DTC2
B	683	SER	ARG	conflict	UNP P0DTC2
B	685	SER	ARG	conflict	UNP P0DTC2
B	817	PRO	PHE	conflict	UNP P0DTC2
B	892	PRO	ALA	conflict	UNP P0DTC2
B	899	PRO	ALA	conflict	UNP P0DTC2
B	942	PRO	ALA	conflict	UNP P0DTC2
B	986	PRO	LYS	conflict	UNP P0DTC2
B	987	PRO	VAL	conflict	UNP P0DTC2
B	1212	SER	-	expression tag	UNP P0DTC2
B	1213	GLY	-	expression tag	UNP P0DTC2
B	1214	GLY	-	expression tag	UNP P0DTC2
B	1215	TYR	-	expression tag	UNP P0DTC2
B	1216	ILE	-	expression tag	UNP P0DTC2
B	1217	PRO	-	expression tag	UNP P0DTC2
B	1218	GLU	-	expression tag	UNP P0DTC2
B	1219	ALA	-	expression tag	UNP P0DTC2
B	1220	PRO	-	expression tag	UNP P0DTC2
B	1221	ARG	-	expression tag	UNP P0DTC2
B	1222	ASP	-	expression tag	UNP P0DTC2
B	1223	GLY	-	expression tag	UNP P0DTC2
B	1224	GLN	-	expression tag	UNP P0DTC2
B	1225	ALA	-	expression tag	UNP P0DTC2
B	1226	TYR	-	expression tag	UNP P0DTC2
B	1227	VAL	-	expression tag	UNP P0DTC2
B	1228	ARG	-	expression tag	UNP P0DTC2
B	1229	LYS	-	expression tag	UNP P0DTC2
B	1230	ASP	-	expression tag	UNP P0DTC2
B	1231	GLY	-	expression tag	UNP P0DTC2
B	1232	GLU	-	expression tag	UNP P0DTC2
B	1233	TRP	-	expression tag	UNP P0DTC2
B	1234	VAL	-	expression tag	UNP P0DTC2
B	1235	LEU	-	expression tag	UNP P0DTC2
B	1236	LEU	-	expression tag	UNP P0DTC2
B	1237	SER	-	expression tag	UNP P0DTC2
B	1238	THR	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1239	PHE	-	expression tag	UNP P0DTC2
B	1240	LEU	-	expression tag	UNP P0DTC2
B	1241	ASN	-	expression tag	UNP P0DTC2
B	1242	SER	-	expression tag	UNP P0DTC2
B	1243	GLY	-	expression tag	UNP P0DTC2
B	1244	ARG	-	expression tag	UNP P0DTC2
B	1245	ALA	-	expression tag	UNP P0DTC2
B	1246	HIS	-	expression tag	UNP P0DTC2
B	1247	HIS	-	expression tag	UNP P0DTC2
B	1248	HIS	-	expression tag	UNP P0DTC2
B	1249	HIS	-	expression tag	UNP P0DTC2
B	1250	HIS	-	expression tag	UNP P0DTC2
B	1251	HIS	-	expression tag	UNP P0DTC2
B	1252	GLY	-	expression tag	UNP P0DTC2
B	1253	ALA	-	expression tag	UNP P0DTC2
B	1254	GLY	-	expression tag	UNP P0DTC2
B	1255	GLY	-	expression tag	UNP P0DTC2
B	1256	LEU	-	expression tag	UNP P0DTC2
B	1257	ASN	-	expression tag	UNP P0DTC2
B	1258	ASP	-	expression tag	UNP P0DTC2
B	1259	ILE	-	expression tag	UNP P0DTC2
B	1260	PHE	-	expression tag	UNP P0DTC2
B	1261	GLU	-	expression tag	UNP P0DTC2
B	1262	ALA	-	expression tag	UNP P0DTC2
B	1263	GLN	-	expression tag	UNP P0DTC2
B	1264	LYS	-	expression tag	UNP P0DTC2
B	1265	ILE	-	expression tag	UNP P0DTC2
B	1266	GLU	-	expression tag	UNP P0DTC2
B	1267	TRP	-	expression tag	UNP P0DTC2
B	1268	HIS	-	expression tag	UNP P0DTC2
B	1269	GLU	-	expression tag	UNP P0DTC2
B	1270	ASP	-	expression tag	UNP P0DTC2
B	1271	THR	-	expression tag	UNP P0DTC2
B	1272	ALA	-	expression tag	UNP P0DTC2
B	1273	ALA	-	expression tag	UNP P0DTC2
B	1274	ALA	-	expression tag	UNP P0DTC2
C	682	SER	ARG	conflict	UNP P0DTC2
C	683	SER	ARG	conflict	UNP P0DTC2
C	685	SER	ARG	conflict	UNP P0DTC2
C	817	PRO	PHE	conflict	UNP P0DTC2
C	892	PRO	ALA	conflict	UNP P0DTC2
C	899	PRO	ALA	conflict	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	942	PRO	ALA	conflict	UNP P0DTC2
C	986	PRO	LYS	conflict	UNP P0DTC2
C	987	PRO	VAL	conflict	UNP P0DTC2
C	1212	SER	-	expression tag	UNP P0DTC2
C	1213	GLY	-	expression tag	UNP P0DTC2
C	1214	GLY	-	expression tag	UNP P0DTC2
C	1215	TYR	-	expression tag	UNP P0DTC2
C	1216	ILE	-	expression tag	UNP P0DTC2
C	1217	PRO	-	expression tag	UNP P0DTC2
C	1218	GLU	-	expression tag	UNP P0DTC2
C	1219	ALA	-	expression tag	UNP P0DTC2
C	1220	PRO	-	expression tag	UNP P0DTC2
C	1221	ARG	-	expression tag	UNP P0DTC2
C	1222	ASP	-	expression tag	UNP P0DTC2
C	1223	GLY	-	expression tag	UNP P0DTC2
C	1224	GLN	-	expression tag	UNP P0DTC2
C	1225	ALA	-	expression tag	UNP P0DTC2
C	1226	TYR	-	expression tag	UNP P0DTC2
C	1227	VAL	-	expression tag	UNP P0DTC2
C	1228	ARG	-	expression tag	UNP P0DTC2
C	1229	LYS	-	expression tag	UNP P0DTC2
C	1230	ASP	-	expression tag	UNP P0DTC2
C	1231	GLY	-	expression tag	UNP P0DTC2
C	1232	GLU	-	expression tag	UNP P0DTC2
C	1233	TRP	-	expression tag	UNP P0DTC2
C	1234	VAL	-	expression tag	UNP P0DTC2
C	1235	LEU	-	expression tag	UNP P0DTC2
C	1236	LEU	-	expression tag	UNP P0DTC2
C	1237	SER	-	expression tag	UNP P0DTC2
C	1238	THR	-	expression tag	UNP P0DTC2
C	1239	PHE	-	expression tag	UNP P0DTC2
C	1240	LEU	-	expression tag	UNP P0DTC2
C	1241	ASN	-	expression tag	UNP P0DTC2
C	1242	SER	-	expression tag	UNP P0DTC2
C	1243	GLY	-	expression tag	UNP P0DTC2
C	1244	ARG	-	expression tag	UNP P0DTC2
C	1245	ALA	-	expression tag	UNP P0DTC2
C	1246	HIS	-	expression tag	UNP P0DTC2
C	1247	HIS	-	expression tag	UNP P0DTC2
C	1248	HIS	-	expression tag	UNP P0DTC2
C	1249	HIS	-	expression tag	UNP P0DTC2
C	1250	HIS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1251	HIS	-	expression tag	UNP P0DTC2
C	1252	GLY	-	expression tag	UNP P0DTC2
C	1253	ALA	-	expression tag	UNP P0DTC2
C	1254	GLY	-	expression tag	UNP P0DTC2
C	1255	GLY	-	expression tag	UNP P0DTC2
C	1256	LEU	-	expression tag	UNP P0DTC2
C	1257	ASN	-	expression tag	UNP P0DTC2
C	1258	ASP	-	expression tag	UNP P0DTC2
C	1259	ILE	-	expression tag	UNP P0DTC2
C	1260	PHE	-	expression tag	UNP P0DTC2
C	1261	GLU	-	expression tag	UNP P0DTC2
C	1262	ALA	-	expression tag	UNP P0DTC2
C	1263	GLN	-	expression tag	UNP P0DTC2
C	1264	LYS	-	expression tag	UNP P0DTC2
C	1265	ILE	-	expression tag	UNP P0DTC2
C	1266	GLU	-	expression tag	UNP P0DTC2
C	1267	TRP	-	expression tag	UNP P0DTC2
C	1268	HIS	-	expression tag	UNP P0DTC2
C	1269	GLU	-	expression tag	UNP P0DTC2
C	1270	ASP	-	expression tag	UNP P0DTC2
C	1271	THR	-	expression tag	UNP P0DTC2
C	1272	ALA	-	expression tag	UNP P0DTC2
C	1273	ALA	-	expression tag	UNP P0DTC2
C	1274	ALA	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called Fab 15033-7 light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	L	214	Total	C	N	O	S	0	0
			1641	1024	276	335	6		
2	M	214	Total	C	N	O	S	0	0
			1641	1024	276	335	6		
2	N	214	Total	C	N	O	S	0	0
			1641	1024	276	335	6		

- Molecule 3 is a protein called Fab 15033-7 heavy chain.

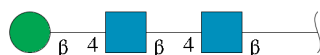
Mol	Chain	Residues	Atoms					AltConf	Trace
3	H	223	Total	C	N	O	S	0	0
			1635	1031	272	324	8		
3	I	223	Total	C	N	O	S	0	0
			1635	1031	272	324	8		

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Mol	Chain	Residues	Atoms					AltConf	Trace
3	J	223	Total	C	N	O	S	0	0
			1635	1031	272	324	8		

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



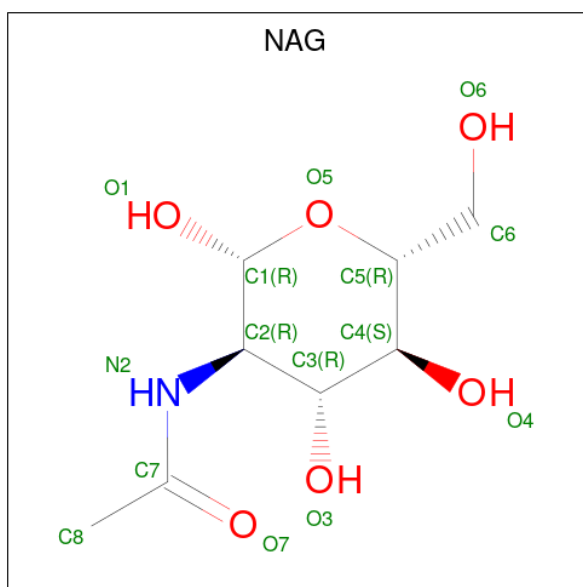
Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	3	Total	C	N	O		0	0
			39	22	2	15			
4	E	3	Total	C	N	O		0	0
			39	22	2	15			

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



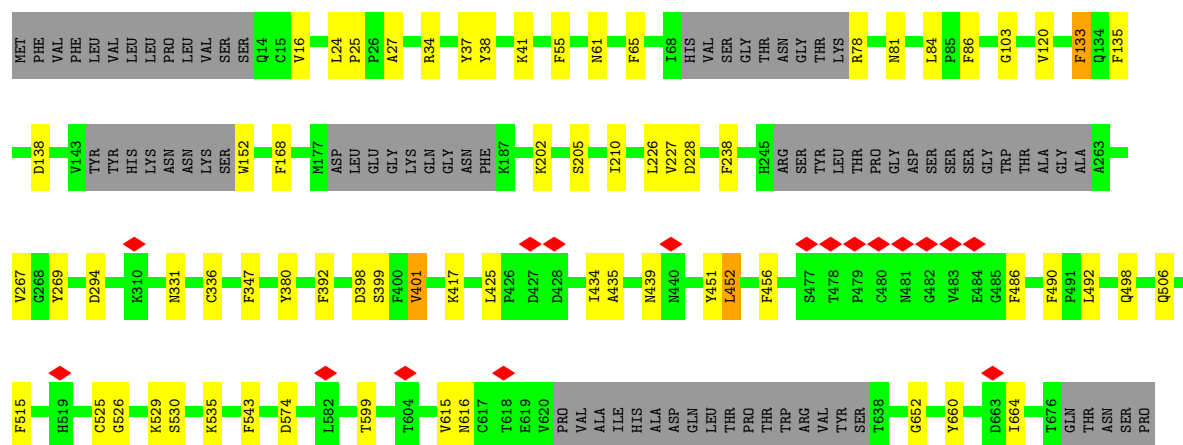
Mol	Chain	Residues	Atoms					AltConf	Trace
5	F	2	Total	C	N	O		0	0
			28	16	2	10			

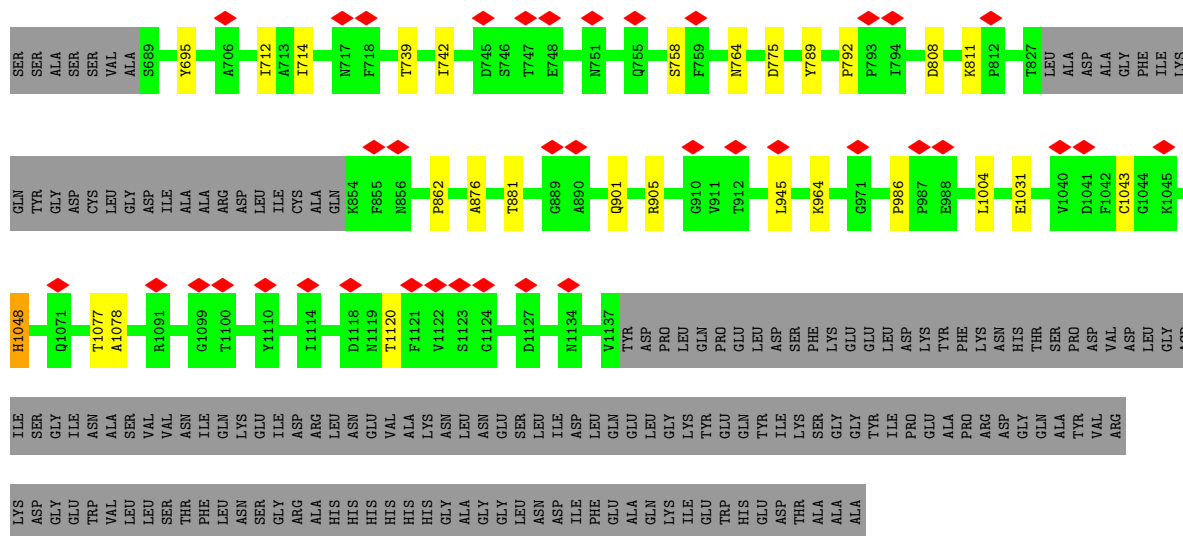
- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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







- Molecule 2: Fab 15033-7 light chain

Chain L: 93% 7%



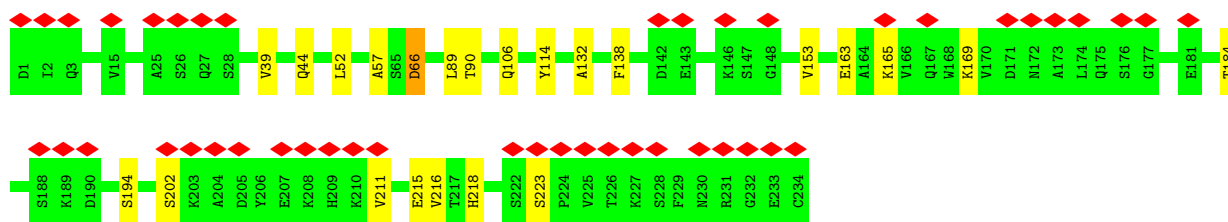
- Molecule 2: Fab 15033-7 light chain

Chain M: 96% 4%



- Molecule 2: Fab 15033-7 light chain

Chain N: 21% 89% 10%

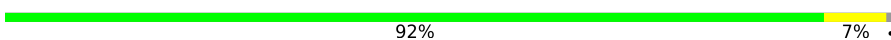


- Molecule 3: Fab 15033-7 heavy chain

Chain H: 93% 5%

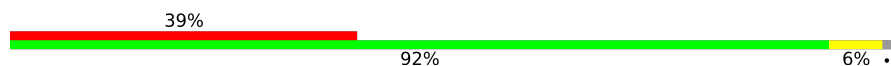


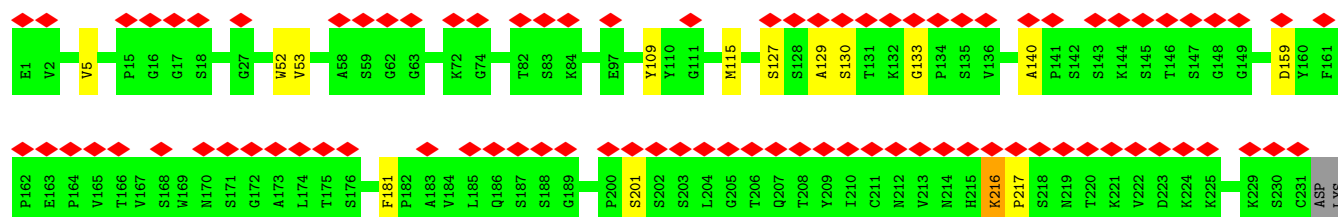
- Molecule 3: Fab 15033-7 heavy chain

Chain I:  92% 7%



- Molecule 3: Fab 15033-7 heavy chain

Chain J:  39% 92% 6%



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

Chain D:  67% 33%



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

Chain E:  33% 67%



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

Chain F:  50% 50%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	19743	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 ($e^-/\text{\AA}^2$)	38.08	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	75000	Depositor
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	1.271	Depositor
Minimum map value	-0.585	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.056	Depositor
Recommended contour level	0.24	Depositor
Map size (Å)	412.0, 412.0, 412.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.03, 1.03, 1.03	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, 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	1.03	6/8200 (0.1%)	1.37	94/11161 (0.8%)
1	B	1.04	5/8200 (0.1%)	1.36	98/11161 (0.9%)
1	C	1.05	7/8200 (0.1%)	1.39	93/11161 (0.8%)
2	L	0.95	0/1676	1.38	21/2276 (0.9%)
2	M	0.99	1/1676 (0.1%)	1.30	9/2276 (0.4%)
2	N	1.01	2/1676 (0.1%)	1.36	15/2276 (0.7%)
3	H	1.00	1/1675 (0.1%)	1.35	16/2280 (0.7%)
3	I	0.98	2/1675 (0.1%)	1.32	17/2280 (0.7%)
3	J	0.99	1/1675 (0.1%)	1.36	16/2280 (0.7%)
All	All	1.03	25/34653 (0.1%)	1.37	379/47151 (0.8%)

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	215	HIS	CB-CG	-8.84	1.37	1.50
3	H	215	HIS	CB-CG	-7.16	1.40	1.50
2	N	89	LEU	CB-CG	-7.08	1.39	1.53
1	C	451	TYR	CB-CG	-6.84	1.36	1.51
1	A	78	ARG	NE-CZ	6.56	1.40	1.33
1	C	78	ARG	NE-CZ	6.34	1.40	1.33
1	B	897	PRO	CA-C	-6.34	1.45	1.52
1	A	801	ASN	C-O	-6.29	1.15	1.23
1	B	152	TRP	CD2-CE3	6.25	1.50	1.40
1	B	152	TRP	CZ2-CH2	6.20	1.49	1.37
1	B	78	ARG	NE-CZ	6.20	1.39	1.33
1	A	152	TRP	CZ2-CH2	6.18	1.49	1.37
1	C	152	TRP	CZ2-CH2	6.15	1.49	1.37
1	C	37	TYR	CB-CG	-6.08	1.38	1.51
1	C	152	TRP	CD2-CE3	6.08	1.50	1.40
1	A	152	TRP	CD2-CE3	6.02	1.49	1.40
2	M	55	TYR	CB-CG	-5.99	1.38	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	109	TYR	CB-CG	-5.95	1.38	1.51
1	A	350	VAL	CB-CG2	-5.78	1.33	1.52
1	B	1048	HIS	CB-CG	-5.45	1.42	1.50
2	N	106	GLN	CG-CD	5.26	1.65	1.52
1	C	1004	LEU	CB-CG	5.17	1.63	1.53
3	I	12	LEU	CB-CG	5.17	1.63	1.53
1	C	78	ARG	CD-NE	5.12	1.53	1.46
1	A	37	TYR	CB-CG	-5.10	1.40	1.51

All (379) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	218	HIS	CA-CB-CG	9.68	123.48	113.80
1	C	486	PHE	CA-CB-CG	-9.29	104.51	113.80
1	A	560	LEU	CA-C-N	8.34	128.00	119.82
1	A	560	LEU	C-N-CA	8.34	128.00	119.82
1	C	16	VAL	N-CA-C	8.26	118.35	110.42
1	C	664	ILE	CA-C-N	8.23	128.23	119.76
1	C	664	ILE	C-N-CA	8.23	128.23	119.76
3	J	130	SER	N-CA-C	8.06	120.97	111.71
1	A	238	PHE	N-CA-C	8.02	121.06	108.79
3	H	137	PHE	CA-C-N	7.93	127.88	120.03
3	H	137	PHE	C-N-CA	7.93	127.88	120.03
1	A	329	PHE	CA-C-N	7.77	128.82	120.11
1	A	329	PHE	C-N-CA	7.77	128.82	120.11
1	C	138	ASP	CA-C-N	7.75	127.74	119.76
1	C	138	ASP	C-N-CA	7.75	127.74	119.76
2	N	90	THR	N-CA-C	7.74	120.83	109.07
1	A	399	SER	N-CA-C	7.66	120.90	108.41
1	B	86	PHE	CA-CB-CG	7.63	121.43	113.80
1	C	210	ILE	N-CA-C	7.62	118.77	108.11
1	B	599	THR	CA-C-N	7.59	128.04	120.52
1	B	599	THR	C-N-CA	7.59	128.04	120.52
1	C	24	LEU	CA-C-N	7.54	125.17	119.66
1	C	24	LEU	C-N-CA	7.54	125.17	119.66
1	C	25	PRO	CA-C-N	7.53	127.70	119.87
1	C	25	PRO	C-N-CA	7.53	127.70	119.87
3	I	137	PHE	CA-C-N	7.53	127.48	120.03
3	I	137	PHE	C-N-CA	7.53	127.48	120.03
1	C	574	ASP	CA-CB-CG	7.51	120.11	112.60
3	I	175	THR	N-CA-C	-7.49	105.05	114.56
1	C	336	CYS	CA-C-N	7.38	127.06	119.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	336	CYS	C-N-CA	7.38	127.06	119.82
1	B	138	ASP	CA-C-N	7.37	127.13	119.76
1	B	138	ASP	C-N-CA	7.37	127.13	119.76
3	J	133	GLY	CA-C-N	7.33	127.49	120.31
3	J	133	GLY	C-N-CA	7.33	127.49	120.31
1	C	1048	HIS	CA-CB-CG	-7.32	106.48	113.80
3	H	6	GLU	N-CA-C	7.30	120.80	110.50
1	B	600	PRO	CA-C-N	-7.28	112.89	120.31
1	B	600	PRO	C-N-CA	-7.28	112.89	120.31
1	A	792	PRO	CA-C-N	7.23	126.86	119.56
1	A	792	PRO	C-N-CA	7.23	126.86	119.56
3	J	53	VAL	N-CA-C	7.19	117.74	111.56
1	A	25	PRO	CA-C-N	7.18	127.34	119.87
1	A	25	PRO	C-N-CA	7.18	127.34	119.87
1	C	425	LEU	CA-C-N	7.09	127.50	119.92
1	C	425	LEU	C-N-CA	7.09	127.50	119.92
1	C	599	THR	CA-C-N	7.01	126.97	120.03
1	C	599	THR	C-N-CA	7.01	126.97	120.03
1	B	329	PHE	CA-C-N	6.98	126.92	120.21
1	B	329	PHE	C-N-CA	6.98	126.92	120.21
1	A	227	VAL	N-CA-C	6.97	117.50	107.88
1	B	236	THR	N-CA-C	-6.97	105.70	114.56
2	M	218	HIS	CA-CB-CG	6.95	120.75	113.80
1	A	1031	GLU	N-CA-C	6.94	118.93	111.36
1	A	24	LEU	CA-C-N	6.94	124.73	119.66
1	A	24	LEU	C-N-CA	6.94	124.73	119.66
3	H	133	GLY	CA-C-N	6.91	127.09	120.31
3	H	133	GLY	C-N-CA	6.91	127.09	120.31
1	B	348	ALA	N-CA-C	6.90	118.90	110.41
1	A	599	THR	CA-C-N	6.88	127.33	120.52
1	A	599	THR	C-N-CA	6.88	127.33	120.52
1	A	60	SER	N-CA-C	-6.87	101.67	110.53
1	A	336	CYS	CA-C-N	6.87	126.55	119.82
1	A	336	CYS	C-N-CA	6.87	126.55	119.82
1	A	666	ILE	N-CA-C	-6.85	106.35	111.90
1	C	543	PHE	CA-CB-CG	6.84	120.64	113.80
1	B	520	ALA	CA-C-N	6.83	126.80	120.03
1	B	520	ALA	C-N-CA	6.83	126.80	120.03
1	C	811	LYS	CA-C-N	6.76	126.46	119.56
1	C	811	LYS	C-N-CA	6.76	126.46	119.56
1	B	483	VAL	N-CA-C	6.76	117.61	107.75
1	B	601	GLY	N-CA-C	-6.71	102.41	111.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	205	SER	N-CA-C	6.70	118.39	108.60
1	C	238	PHE	N-CA-C	6.70	119.26	109.07
1	A	1048	HIS	CA-CB-CG	-6.68	107.12	113.80
1	B	811	LYS	CA-C-N	6.64	126.34	119.56
1	B	811	LYS	C-N-CA	6.64	126.34	119.56
2	N	218	HIS	CA-CB-CG	6.63	120.43	113.80
1	C	55	PHE	CA-CB-CG	6.59	120.39	113.80
1	B	792	PRO	CA-C-N	6.58	126.21	119.56
1	B	792	PRO	C-N-CA	6.58	126.21	119.56
1	A	490	PHE	N-CA-C	-6.58	102.03	109.60
1	B	159	VAL	N-CA-C	-6.57	105.91	111.56
1	A	811	LYS	CA-C-N	6.56	126.25	119.56
1	A	811	LYS	C-N-CA	6.56	126.25	119.56
1	B	717	ASN	CA-CB-CG	6.56	119.16	112.60
1	C	267	VAL	N-CA-C	6.55	118.15	108.45
1	A	456	PHE	CA-CB-CG	6.50	120.31	113.80
1	A	945	LEU	CA-C-N	6.50	128.27	120.14
1	A	945	LEU	C-N-CA	6.50	128.27	120.14
1	C	1078	ALA	CA-C-N	6.49	126.12	119.56
1	C	1078	ALA	C-N-CA	6.49	126.12	119.56
1	C	401	VAL	N-CA-C	6.48	117.18	108.11
2	N	132	ALA	CA-C-N	6.48	126.39	119.85
2	N	132	ALA	C-N-CA	6.48	126.39	119.85
1	B	802	PHE	N-CA-C	-6.41	105.22	113.16
1	C	133	PHE	CA-CB-CG	6.39	120.19	113.80
1	A	84	LEU	CA-C-N	6.39	126.36	119.78
1	A	84	LEU	C-N-CA	6.39	126.36	119.78
2	N	216	VAL	N-CA-C	6.37	117.04	107.80
1	C	84	LEU	CA-C-N	6.33	126.30	119.78
1	C	84	LEU	C-N-CA	6.33	126.30	119.78
1	B	498	GLN	CA-C-N	6.31	126.44	119.87
1	B	498	GLN	C-N-CA	6.31	126.44	119.87
2	L	139	PRO	N-CA-C	-6.31	104.41	110.47
1	A	159	VAL	N-CA-C	-6.30	105.59	111.45
1	A	267	VAL	N-CA-C	6.30	117.19	108.12
1	A	737	ASP	CA-C-N	6.26	128.99	120.54
1	A	737	ASP	C-N-CA	6.26	128.99	120.54
1	B	55	PHE	CA-CB-CG	6.26	120.06	113.80
3	H	159	ASP	N-CA-C	6.22	119.97	111.39
1	B	1048	HIS	CA-CB-CG	-6.21	107.59	113.80
1	B	862	PRO	CA-C-N	6.20	126.39	120.31
1	B	862	PRO	C-N-CA	6.20	126.39	120.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	435	ALA	N-CA-C	6.19	119.00	108.90
1	B	800	PHE	CA-CB-CG	6.18	119.98	113.80
1	B	1078	ALA	CA-C-N	6.18	125.80	119.56
1	B	1078	ALA	C-N-CA	6.18	125.80	119.56
1	C	227	VAL	N-CA-C	6.18	116.41	107.88
2	M	202	SER	N-CA-C	-6.18	102.81	110.41
1	C	792	PRO	CA-C-N	6.18	125.80	119.56
1	C	792	PRO	C-N-CA	6.18	125.80	119.56
1	C	399	SER	N-CA-C	6.15	118.93	108.90
1	B	560	LEU	CA-C-N	6.14	126.94	120.12
1	B	560	LEU	C-N-CA	6.14	126.94	120.12
1	B	571	ASP	N-CA-C	6.14	120.03	111.74
2	N	202	SER	N-CA-C	-6.13	102.27	110.55
1	B	576	VAL	N-CA-C	6.13	116.95	108.12
3	H	130	SER	N-CA-C	6.12	120.01	112.54
1	C	120	VAL	N-CA-C	6.12	116.93	108.12
2	L	132	ALA	CA-C-N	6.12	126.03	119.85
2	L	132	ALA	C-N-CA	6.12	126.03	119.85
1	B	714	ILE	CA-C-N	6.10	126.05	119.76
1	B	714	ILE	C-N-CA	6.10	126.05	119.76
1	C	81	ASN	CA-C-N	6.10	126.07	120.03
1	C	81	ASN	C-N-CA	6.10	126.07	120.03
1	A	120	VAL	N-CA-C	6.09	116.89	108.12
1	C	86	PHE	CA-CB-CG	6.08	119.88	113.80
1	C	808	ASP	CA-C-N	6.07	126.18	119.87
1	C	808	ASP	C-N-CA	6.07	126.18	119.87
1	C	498	GLN	CA-C-N	6.06	126.18	119.87
1	C	498	GLN	C-N-CA	6.06	126.18	119.87
1	B	409	GLN	CA-C-N	-6.06	114.90	122.48
1	B	409	GLN	C-N-CA	-6.06	114.90	122.48
2	M	138	PHE	CA-C-N	6.06	124.08	119.66
2	M	138	PHE	C-N-CA	6.06	124.08	119.66
1	C	294	ASP	CA-C-N	6.06	125.94	119.28
1	C	294	ASP	C-N-CA	6.06	125.94	119.28
1	B	383	SER	N-CA-C	-6.05	102.14	109.83
3	J	216	LYS	CA-C-N	6.05	125.73	119.56
3	J	216	LYS	C-N-CA	6.05	125.73	119.56
1	A	86	PHE	CA-CB-CG	6.03	119.83	113.80
1	C	526	GLY	CA-C-N	6.02	125.90	119.28
1	C	526	GLY	C-N-CA	6.02	125.90	119.28
1	A	94	SER	N-CA-C	6.01	118.29	108.55
1	B	808	ASP	CA-C-N	6.00	126.11	119.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	808	ASP	C-N-CA	6.00	126.11	119.87
1	C	764	ASN	N-CA-C	-5.98	104.76	111.28
1	A	506	GLN	CA-C-N	5.96	125.87	119.85
1	A	506	GLN	C-N-CA	5.96	125.87	119.85
1	A	402	ILE	N-CA-C	5.96	117.49	108.80
1	B	883	THR	N-CA-C	5.93	121.30	112.94
2	L	138	PHE	CA-C-N	5.88	123.95	119.66
2	L	138	PHE	C-N-CA	5.88	123.95	119.66
3	J	201	SER	N-CA-C	5.88	118.47	111.71
2	L	223	SER	CA-C-N	5.87	125.83	119.78
2	L	223	SER	C-N-CA	5.87	125.83	119.78
1	B	888	PHE	CA-CB-CG	5.83	119.64	113.80
1	B	84	LEU	CA-C-N	5.83	125.77	119.76
1	B	84	LEU	C-N-CA	5.83	125.77	119.76
1	B	79	PHE	N-CA-C	5.82	117.70	108.67
3	J	127	SER	N-CA-C	5.82	117.95	108.76
2	N	211	VAL	N-CA-C	5.81	116.49	108.12
3	H	216	LYS	CA-C-N	5.80	125.48	119.56
3	H	216	LYS	C-N-CA	5.80	125.48	119.56
1	A	808	ASP	CA-C-N	5.80	125.90	119.87
1	A	808	ASP	C-N-CA	5.80	125.90	119.87
1	A	652	GLY	N-CA-C	-5.78	106.48	115.67
1	A	205	SER	N-CA-C	5.77	117.62	108.79
3	I	116	ASP	CA-CB-CG	5.77	118.37	112.60
1	C	652	GLY	N-CA-C	-5.75	106.57	116.01
2	L	95	GLN	CA-C-N	5.75	125.88	119.32
2	L	95	GLN	C-N-CA	5.75	125.88	119.32
1	A	661	GLU	N-CA-C	-5.75	102.39	110.50
1	C	506	GLN	CA-C-N	5.75	125.66	119.85
1	C	506	GLN	C-N-CA	5.75	125.66	119.85
1	B	298	GLU	N-CA-C	-5.75	105.09	111.36
1	B	945	LEU	CA-C-N	5.74	127.32	120.14
1	B	945	LEU	C-N-CA	5.74	127.32	120.14
1	B	1042	PHE	N-CA-C	5.74	117.21	111.07
1	A	900	MET	N-CA-C	-5.74	105.39	112.90
2	N	223	SER	CA-C-N	5.73	125.74	119.90
2	N	223	SER	C-N-CA	5.73	125.74	119.90
3	J	159	ASP	N-CA-C	5.72	119.29	111.39
1	C	739	THR	N-CA-C	-5.71	105.42	112.90
1	A	138	ASP	CA-CB-CG	5.71	118.31	112.60
1	C	1077	THR	N-CA-C	5.71	118.08	109.41
2	L	233	GLU	CA-C-N	5.71	131.97	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	233	GLU	C-N-CA	5.71	131.97	121.70
1	C	34	ARG	N-CA-C	5.70	119.29	111.54
1	A	1077	THR	N-CA-C	5.69	118.05	109.41
1	A	117	LEU	N-CA-C	5.68	117.99	108.73
1	B	338	PHE	CA-CB-CG	5.67	119.47	113.80
1	C	742	ILE	N-CA-C	5.65	115.84	110.53
1	B	220	PHE	CA-CB-CG	-5.64	108.16	113.80
1	A	748	GLU	CA-C-N	5.63	127.77	120.44
1	A	748	GLU	C-N-CA	5.63	127.77	120.44
1	A	208	THR	CA-C-N	5.63	125.58	119.78
1	A	208	THR	C-N-CA	5.63	125.58	119.78
1	A	393	THR	N-CA-C	-5.63	105.06	111.14
2	N	66	ASP	N-CA-C	5.61	118.62	108.69
1	A	434	ILE	N-CA-CB	-5.61	104.61	111.67
1	B	1068	VAL	CA-C-N	5.61	125.93	119.93
1	B	1068	VAL	C-N-CA	5.61	125.93	119.93
1	B	81	ASN	CA-C-N	5.60	125.58	120.03
1	B	81	ASN	C-N-CA	5.60	125.58	120.03
2	L	202	SER	N-CA-C	-5.60	103.14	110.53
1	C	38	TYR	CA-C-N	5.59	125.43	119.28
1	C	38	TYR	C-N-CA	5.59	125.43	119.28
1	A	526	GLY	N-CA-C	-5.58	100.95	112.34
2	M	132	ALA	CA-C-N	5.58	125.48	119.85
2	M	132	ALA	C-N-CA	5.58	125.48	119.85
1	B	38	TYR	CA-C-N	5.58	125.67	119.87
1	B	38	TYR	C-N-CA	5.58	125.67	119.87
1	B	267	VAL	N-CA-C	5.58	116.31	108.17
1	C	439	ASN	CA-CB-CG	5.58	118.17	112.60
1	A	1057	PRO	N-CA-C	-5.57	102.44	111.19
1	A	742	ILE	N-CA-C	5.57	117.07	111.00
2	L	134	SER	CA-C-N	-5.57	116.29	123.19
2	L	134	SER	C-N-CA	-5.57	116.29	123.19
2	N	44	GLN	N-CA-C	5.56	117.97	108.90
1	C	945	LEU	CA-C-N	5.56	127.09	120.14
1	C	945	LEU	C-N-CA	5.56	127.09	120.14
1	C	331	ASN	CA-CB-CG	5.55	118.16	112.60
3	H	200	PRO	N-CA-C	-5.55	102.38	111.15
3	I	133	GLY	CA-C-N	5.55	125.75	120.31
3	I	133	GLY	C-N-CA	5.55	125.75	120.31
1	A	216	LEU	CA-C-N	5.55	125.55	119.89
1	A	216	LEU	C-N-CA	5.55	125.55	119.89
1	B	452	LEU	N-CA-C	5.54	118.61	109.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	342	PHE	N-CA-C	-5.54	102.69	110.50
3	H	227	GLU	CA-C-N	5.53	125.48	119.78
3	H	227	GLU	C-N-CA	5.53	125.48	119.78
3	J	140	ALA	CA-C-N	5.53	125.50	120.03
3	J	140	ALA	C-N-CA	5.53	125.50	120.03
1	B	173	GLN	CA-C-N	5.52	125.42	119.85
1	B	173	GLN	C-N-CA	5.52	125.42	119.85
1	C	84	LEU	CB-CA-C	5.50	117.14	108.88
2	N	89	LEU	N-CA-CB	-5.50	101.91	110.55
1	C	226	LEU	N-CA-C	-5.50	105.62	112.93
2	L	139	PRO	CA-C-O	-5.49	116.95	120.90
1	B	1031	GLU	N-CA-C	5.48	119.14	112.23
1	C	615	VAL	CA-C-N	5.48	128.65	120.87
1	C	615	VAL	C-N-CA	5.48	128.65	120.87
3	J	5	VAL	N-CA-C	5.47	115.74	107.75
1	C	714	ILE	CA-C-N	5.47	125.39	119.76
1	C	714	ILE	C-N-CA	5.47	125.39	119.76
1	B	205	SER	N-CA-C	5.46	117.38	109.07
2	L	162	ARG	N-CA-C	5.46	117.98	111.71
1	C	490	PHE	CA-C-N	5.45	125.07	119.56
1	C	490	PHE	C-N-CA	5.45	125.07	119.56
1	C	862	PRO	CA-C-N	5.45	125.92	120.14
1	C	862	PRO	C-N-CA	5.45	125.92	120.14
1	A	498	GLN	CA-C-N	5.44	125.53	119.87
1	A	498	GLN	C-N-CA	5.44	125.53	119.87
2	N	138	PHE	CA-C-N	5.44	125.98	120.38
2	N	138	PHE	C-N-CA	5.44	125.98	120.38
1	A	777	ASN	CA-CB-CG	5.42	118.02	112.60
1	A	1052	PHE	CA-C-N	5.41	125.87	120.14
1	A	1052	PHE	C-N-CA	5.41	125.87	120.14
1	C	758	SER	N-CA-C	5.40	118.88	112.72
3	I	130	SER	N-CA-C	5.40	119.13	112.54
1	B	581	THR	N-CA-C	5.40	117.62	109.41
1	A	90	VAL	N-CA-C	5.40	117.86	108.90
1	B	801	ASN	CA-CB-CG	5.40	118.00	112.60
1	B	326	ILE	N-CA-C	5.38	115.60	107.75
1	B	215	ASP	CA-CB-CG	5.37	117.97	112.60
1	C	901	GLN	N-CA-C	-5.37	105.51	111.36
1	A	229	LEU	CA-C-N	5.36	126.54	119.84
1	A	229	LEU	C-N-CA	5.36	126.54	119.84
1	A	231	ILE	N-CA-C	5.36	116.76	111.67
1	A	236	THR	N-CA-C	-5.34	106.64	114.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	269	TYR	N-CA-C	5.34	118.27	109.72
1	A	361	CYS	N-CA-C	5.33	115.58	108.38
1	A	168	PHE	N-CA-C	5.32	117.85	108.75
3	J	181	PHE	CA-C-N	5.32	125.22	119.85
3	J	181	PHE	C-N-CA	5.32	125.22	119.85
1	B	1074	ASN	CA-CB-CG	5.30	117.90	112.60
1	C	775	ASP	CA-CB-CG	5.30	117.90	112.60
1	C	712	ILE	N-CA-C	5.29	115.52	108.11
1	B	329	PHE	CA-CB-CG	5.29	119.09	113.80
1	B	861	LEU	CA-C-N	5.27	123.45	119.66
1	B	861	LEU	C-N-CA	5.27	123.45	119.66
1	B	727	LEU	CA-C-N	5.27	125.03	119.76
1	B	727	LEU	C-N-CA	5.27	125.03	119.76
1	B	1077	THR	N-CA-C	5.26	117.81	109.50
1	B	208	THR	CA-C-N	5.24	125.18	119.78
1	B	208	THR	C-N-CA	5.24	125.18	119.78
1	C	529	LYS	CA-C-N	5.22	129.65	121.08
1	C	529	LYS	C-N-CA	5.22	129.65	121.08
3	H	83	SER	N-CA-C	-5.21	106.07	112.90
1	A	65	PHE	CA-C-N	5.21	128.26	120.87
1	A	65	PHE	C-N-CA	5.21	128.26	120.87
1	A	329	PHE	CA-CB-CG	-5.21	108.59	113.80
1	B	806	LEU	CA-C-N	5.21	125.11	119.85
1	B	806	LEU	C-N-CA	5.21	125.11	119.85
1	B	399	SER	N-CA-C	5.21	117.97	109.59
1	A	271	GLN	CA-C-N	5.20	126.34	119.84
1	A	271	GLN	C-N-CA	5.20	126.34	119.84
1	A	712	ILE	N-CA-C	5.20	115.39	108.11
3	I	216	LYS	CA-C-N	5.20	124.87	119.56
3	I	216	LYS	C-N-CA	5.20	124.87	119.56
1	C	434	ILE	N-CA-CB	-5.19	104.90	111.64
2	M	139	PRO	N-CA-C	-5.19	105.49	110.47
1	C	1031	GLU	N-CA-C	5.18	119.69	112.90
2	L	29	VAL	N-CA-C	-5.18	106.39	111.88
1	A	103	GLY	N-CA-C	5.18	117.58	110.69
1	B	709	ASN	CA-CB-CG	5.18	117.78	112.60
1	A	497	PHE	CA-CB-CG	5.17	118.97	113.80
1	A	81	ASN	CA-C-N	5.17	125.15	120.03
1	A	81	ASN	C-N-CA	5.17	125.15	120.03
1	B	133	PHE	CA-CB-CG	5.17	118.97	113.80
1	A	1040	VAL	CA-C-N	5.16	129.67	122.08
1	A	1040	VAL	C-N-CA	5.16	129.67	122.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	530	SER	N-CA-C	-5.16	105.27	112.03
1	B	425	LEU	CA-C-N	5.15	125.09	119.78
1	B	425	LEU	C-N-CA	5.15	125.09	119.78
1	C	103	GLY	N-CA-C	5.14	118.64	111.19
1	C	1120	THR	N-CA-C	5.13	117.47	109.52
3	I	83	SER	N-CA-C	-5.13	105.76	112.23
3	I	88	TYR	N-CA-C	5.13	117.61	109.50
1	A	1042	PHE	N-CA-C	5.13	116.56	111.07
2	L	92	SER	N-CA-C	5.13	117.77	111.82
3	I	159	ASP	N-CA-C	5.12	118.66	111.74
1	A	374	PHE	CA-CB-CG	-5.12	108.68	113.80
1	B	495	TYR	CB-CA-C	-5.12	101.31	109.75
3	I	181	PHE	CA-C-N	5.12	125.02	119.85
3	I	181	PHE	C-N-CA	5.12	125.02	119.85
1	B	238	PHE	N-CA-C	5.11	116.84	109.07
1	A	562	PHE	CA-CB-CG	5.11	118.91	113.80
1	B	1039	ARG	N-CA-C	-5.11	102.30	109.96
1	B	664	ILE	CA-C-N	5.11	125.02	119.76
1	B	664	ILE	C-N-CA	5.11	125.02	119.76
2	N	153	VAL	N-CA-C	5.11	115.26	108.11
1	C	168	PHE	N-CA-C	5.10	117.47	108.75
2	L	153	VAL	N-CA-C	5.09	115.24	108.11
1	B	155	SER	N-CA-C	5.09	117.64	108.48
1	C	452	LEU	CA-C-N	5.09	128.78	121.50
1	C	452	LEU	C-N-CA	5.09	128.78	121.50
1	B	495	TYR	CA-CB-CG	5.09	123.06	113.90
1	C	398	ASP	N-CA-CB	-5.09	102.38	111.08
1	C	456	PHE	N-CA-C	5.09	117.26	109.07
1	B	110	LEU	CB-CA-C	-5.08	103.08	111.36
3	J	129	ALA	CA-C-N	5.08	127.59	120.28
3	J	129	ALA	C-N-CA	5.08	127.59	120.28
3	H	140	ALA	CA-C-N	5.07	125.05	120.03
3	H	140	ALA	C-N-CA	5.07	125.05	120.03
3	H	5	VAL	N-CA-C	5.07	115.21	108.11
1	A	157	PHE	CA-CB-CG	-5.07	108.73	113.80
1	A	237	ARG	N-CA-C	5.07	116.89	109.24
1	A	670	ILE	CA-C-N	5.07	131.96	122.74
1	A	670	ILE	C-N-CA	5.07	131.96	122.74
1	C	986	PRO	N-CA-C	5.06	116.88	110.70
2	M	216	VAL	N-CA-C	5.05	115.37	107.99
2	M	44	GLN	N-CA-C	5.05	117.13	108.90
3	I	106	ARG	N-CA-C	5.04	116.86	109.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	338	PHE	N-CA-C	-5.04	106.23	112.38
1	A	985	ASP	CA-C-N	5.04	125.57	120.38
1	A	985	ASP	C-N-CA	5.04	125.57	120.38
1	B	216	LEU	CA-C-N	5.04	124.94	119.85
1	B	216	LEU	C-N-CA	5.04	124.94	119.85
1	A	567	ARG	N-CA-C	5.03	117.45	109.50
1	B	892	PRO	N-CA-C	-5.03	103.21	111.21
1	C	65	PHE	CA-C-N	5.02	128.00	120.87
1	C	65	PHE	C-N-CA	5.02	128.00	120.87
2	L	134	SER	N-CA-C	-5.02	101.57	109.25
1	B	1062	PHE	CA-CB-CG	5.01	118.81	113.80
3	I	227	GLU	CA-C-N	5.00	124.91	119.76
3	I	227	GLU	C-N-CA	5.00	124.91	119.76

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	8017	0	7835	14	0
1	B	8017	0	7834	16	0
1	C	8017	0	7835	16	0
2	L	1641	0	1594	2	0
2	M	1641	0	1594	0	0
2	N	1641	0	1594	8	0
3	H	1635	0	1586	2	0
3	I	1635	0	1586	1	0
3	J	1635	0	1586	3	0
4	D	39	0	34	0	0
4	E	39	0	34	0	0
5	F	28	0	25	0	0
6	A	28	0	26	0	0
6	B	28	0	26	0	0
6	C	42	0	39	2	0
All	All	34083	0	33228	54	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 1.

All (54) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:27:ALA:HB1	6:C:5002:NAG:H2	1.79	0.65
1:C:61:ASN:O	6:C:5002:NAG:H82	2.01	0.60
1:C:417:LYS:NZ	2:N:66:ASP:OD1	2.37	0.57
2:L:142:ASP:OD1	2:L:142:ASP:N	2.39	0.55
1:C:202:LYS:NZ	1:C:228:ASP:OD1	2.41	0.52
1:B:811:LYS:NZ	1:B:820:ASP:OD2	2.43	0.52
1:A:854:LYS:NZ	1:B:568:ASP:OD2	2.43	0.51
1:C:452:LEU:HB3	1:C:492:LEU:HB3	1.91	0.51
3:H:6:GLU:H	3:H:6:GLU:CD	2.18	0.51
1:A:571:ASP:OD2	1:C:964:LYS:NZ	2.40	0.51
1:A:977:LEU:HD11	1:A:993:ILE:HG12	1.92	0.51
1:A:489:TYR:O	1:A:490:PHE:C	2.52	0.50
2:N:163:GLU:OE2	2:N:165:LYS:NZ	2.45	0.50
1:A:43:PHE:CE2	1:B:563:GLN:HB2	2.47	0.49
2:N:39:VAL:H	2:N:57:ALA:H	1.58	0.49
1:A:202:LYS:NZ	1:A:228:ASP:OD1	2.47	0.48
2:L:165:LYS:NZ	2:L:181:GLU:OE2	2.48	0.47
3:I:216:LYS:N	3:I:217:PRO:CD	2.78	0.47
1:A:30:ASN:OD1	1:A:30:ASN:C	2.56	0.47
1:C:347:PHE:CD1	1:C:401:VAL:HB	2.51	0.46
2:N:169:LYS:NZ	2:N:215:GLU:OE1	2.49	0.46
1:C:133:PHE:HB2	1:C:135:PHE:CE1	2.51	0.46
1:B:349:SER:HB3	1:B:351:TYR:CE2	2.51	0.45
3:H:216:LYS:N	3:H:217:PRO:CD	2.79	0.45
1:A:563:GLN:HB3	1:C:41:LYS:O	2.15	0.45
1:B:55:PHE:HB2	1:B:275:PHE:CE1	2.52	0.45
1:A:964:LYS:NZ	1:B:571:ASP:OD2	2.50	0.45
1:C:347:PHE:CE1	1:C:401:VAL:HB	2.51	0.45
1:B:57:PRO:HA	1:B:290:ASP:OD2	2.18	0.44
1:A:725:GLU:OE1	1:A:1028:LYS:NZ	2.50	0.44
1:A:1043:CYS:H	1:A:1048:HIS:CE1	2.35	0.44
1:B:329:PHE:CG	1:B:330:PRO:HD2	2.53	0.44
1:B:389:ASP:N	1:B:389:ASP:OD1	2.48	0.43
1:B:1043:CYS:HB2	1:B:1048:HIS:CE1	2.53	0.43
3:J:216:LYS:N	3:J:217:PRO:CD	2.81	0.43
1:C:660:TYR:CD1	1:C:695:TYR:CD2	3.05	0.43
2:N:184:THR:HG22	2:N:194:SER:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1043:CYS:H	1:C:1048:HIS:CE1	2.38	0.42
1:A:61:ASN:OD1	1:A:61:ASN:N	2.53	0.41
1:B:516:GLU:HG3	1:B:518:LEU:H	1.84	0.41
1:B:989:ALA:O	1:B:993:ILE:HG12	2.20	0.41
1:C:881:THR:O	1:C:905:ARG:NH2	2.53	0.41
2:N:52:LEU:HD22	3:J:115:MET:O	2.21	0.41
1:B:400:PHE:CD2	1:B:400:PHE:N	2.88	0.41
1:B:568:ASP:OD1	1:B:568:ASP:C	2.62	0.41
2:N:114:TYR:CD2	3:J:52:TRP:CH2	3.08	0.41
1:C:380:TYR:CD1	1:C:380:TYR:N	2.88	0.41
1:C:789:TYR:H	1:C:876:ALA:CB	2.33	0.41
2:N:184:THR:CG2	2:N:194:SER:H	2.34	0.41
1:A:438:SER:HB3	1:A:509:ARG:HG2	2.02	0.41
1:A:733:LYS:NZ	1:A:775:ASP:OD1	2.34	0.40
1:C:392:PHE:CD2	1:C:515:PHE:HB3	2.56	0.40
1:B:351:TYR:CD2	1:B:351:TYR:N	2.89	0.40
1:B:449:TYR:CD2	1:B:494:SER:HB3	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	1010/1274 (79%)	998 (99%)	12 (1%)	0	100	100
1	B	1010/1274 (79%)	990 (98%)	19 (2%)	1 (0%)	48	83
1	C	1010/1274 (79%)	992 (98%)	16 (2%)	2 (0%)	43	77
2	L	212/214 (99%)	210 (99%)	2 (1%)	0	100	100
2	M	212/214 (99%)	209 (99%)	3 (1%)	0	100	100
2	N	212/214 (99%)	211 (100%)	1 (0%)	0	100	100
3	H	221/225 (98%)	216 (98%)	5 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	I	221/225 (98%)	217 (98%)	4 (2%)	0	100	100
3	J	221/225 (98%)	216 (98%)	5 (2%)	0	100	100
All	All	4329/5139 (84%)	4259 (98%)	67 (2%)	3 (0%)	49	83

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	337	PRO
1	C	525	CYS
1	C	535	LYS

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	901/1107 (81%)	901 (100%)	0	100	100
1	B	901/1107 (81%)	897 (100%)	4 (0%)	84	84
1	C	901/1107 (81%)	900 (100%)	1 (0%)	88	89
2	L	189/189 (100%)	189 (100%)	0	100	100
2	M	189/189 (100%)	189 (100%)	0	100	100
2	N	189/189 (100%)	189 (100%)	0	100	100
3	H	177/179 (99%)	177 (100%)	0	100	100
3	I	177/179 (99%)	175 (99%)	2 (1%)	65	76
3	J	177/179 (99%)	177 (100%)	0	100	100
All	All	3801/4425 (86%)	3794 (100%)	7 (0%)	85	87

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	61	ASN
1	B	335	LEU
1	B	657	ASN

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Mol	Chain	Res	Type
1	B	717	ASN
1	C	616	ASN
3	I	23	CYS
3	I	104	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	616	ASN
1	A	657	ASN
1	A	777	ASN
1	A	801	ASN
1	A	920	GLN
1	A	935	GLN
1	B	122	ASN
1	B	282	ASN
1	B	493	GLN
1	B	690	GLN
1	B	717	ASN
1	B	824	ASN
1	B	901	GLN
1	B	920	GLN
1	B	965	GLN
1	B	1083	HIS
1	C	196	ASN
1	C	717	ASN
1	C	913	GLN
1	C	920	GLN
2	L	175	GLN
2	L	178	ASN
3	H	214	ASN
2	M	167	GLN
3	I	214	ASN
3	J	214	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

8 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	D	1	4,1	14,14,15	0.41	0	17,19,21	1.46	2 (11%)
4	NAG	D	2	4	14,14,15	0.36	0	17,19,21	0.79	0
4	BMA	D	3	4	11,11,12	0.30	0	15,15,17	0.55	0
4	NAG	E	1	4,1	14,14,15	0.50	0	17,19,21	2.45	5 (29%)
4	NAG	E	2	4	14,14,15	0.35	0	17,19,21	0.89	2 (11%)
4	BMA	E	3	4	11,11,12	0.28	0	15,15,17	0.50	0
5	NAG	F	1	5,1	14,14,15	0.35	0	17,19,21	1.61	2 (11%)
5	NAG	F	2	5	14,14,15	0.33	0	17,19,21	0.62	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	D	1	4,1	-	4/6/23/26	0/1/1/1
4	NAG	D	2	4	-	0/6/23/26	0/1/1/1
4	BMA	D	3	4	-	2/2/19/22	0/1/1/1
4	NAG	E	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	2/2/19/22	0/1/1/1
5	NAG	F	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	F	2	5	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	1	NAG	C1-O5-C5	5.00	118.89	112.19
4	E	1	NAG	C4-C3-C2	-4.89	103.85	111.02
4	E	1	NAG	C3-C4-C5	-4.85	101.44	110.23
4	E	1	NAG	C1-O5-C5	4.79	118.61	112.19
4	D	1	NAG	C3-C4-C5	3.69	116.93	110.23
4	E	1	NAG	C2-N2-C7	-3.35	118.42	122.90
4	D	1	NAG	O5-C1-C2	-3.34	106.12	111.29
5	F	1	NAG	C4-C3-C2	-3.20	106.32	111.02
4	E	1	NAG	C1-C2-N2	3.02	115.19	110.43
4	E	2	NAG	O5-C1-C2	-2.19	107.90	111.29
4	E	2	NAG	C2-N2-C7	-2.11	120.08	122.90

There are no chirality outliers.

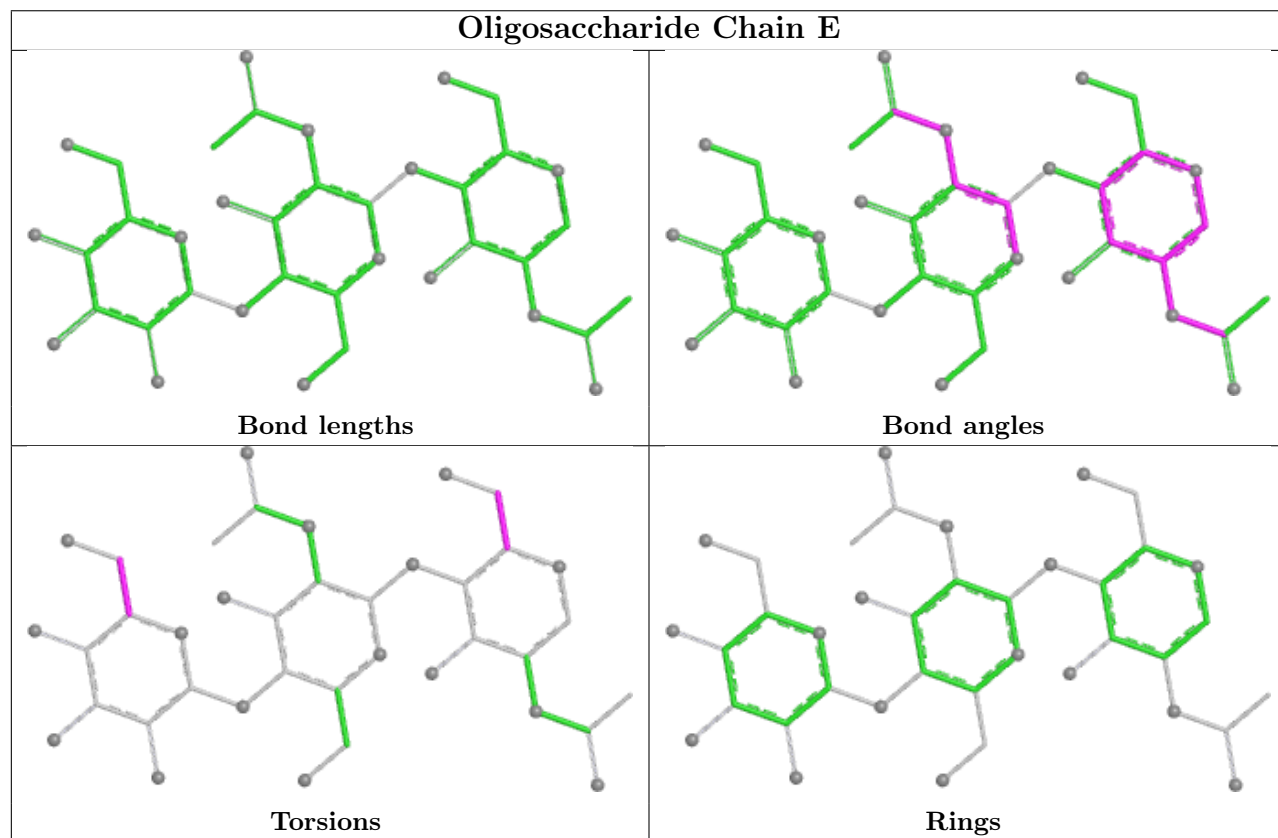
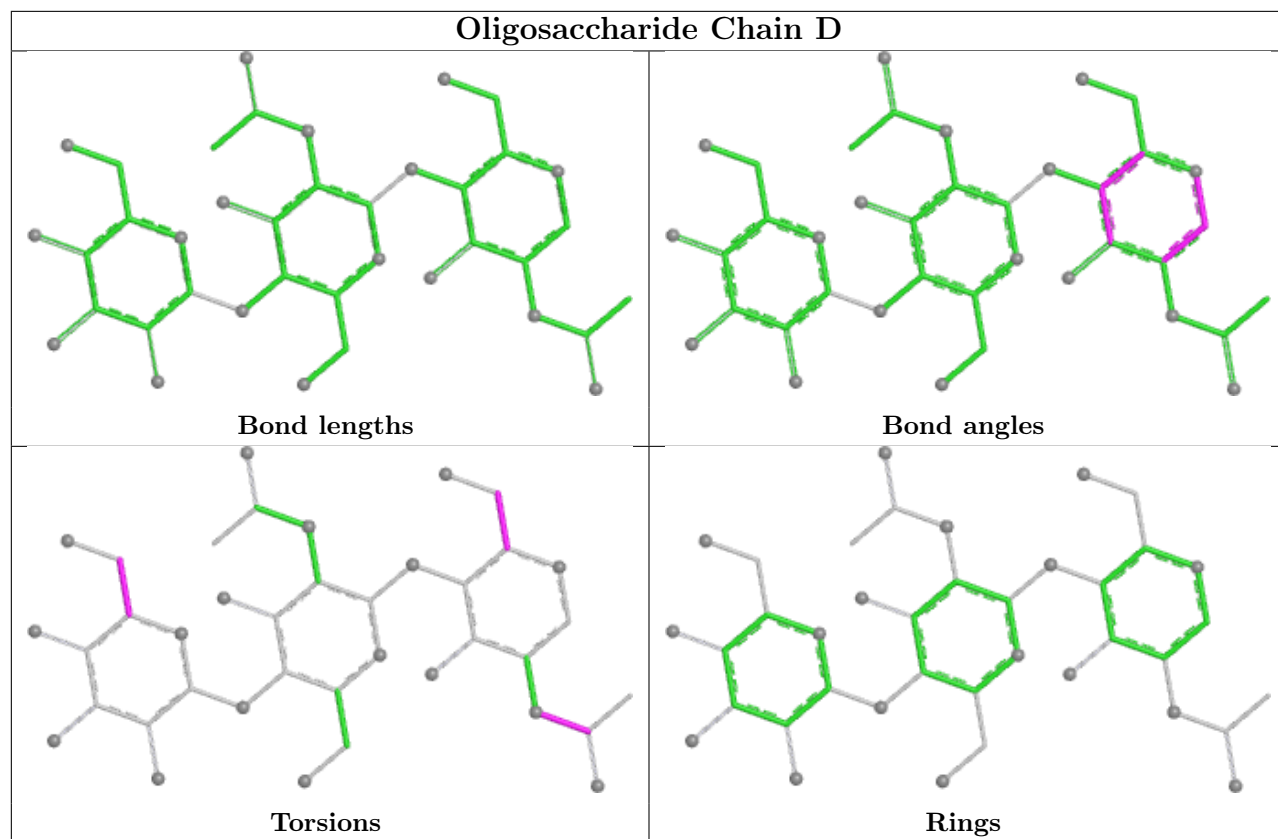
All (14) torsion outliers are listed below:

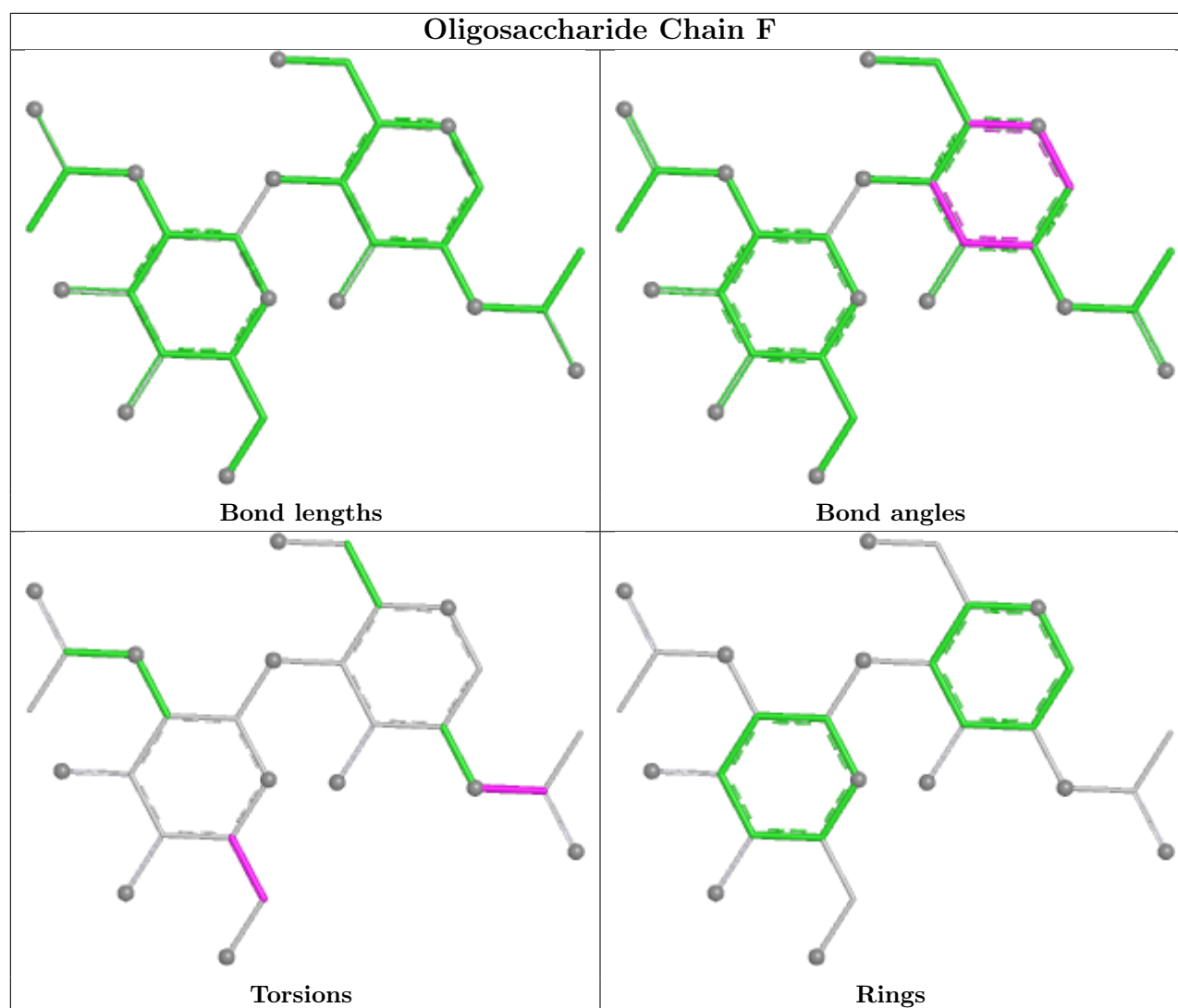
Mol	Chain	Res	Type	Atoms
4	E	3	BMA	O5-C5-C6-O6
4	E	1	NAG	O5-C5-C6-O6
4	D	3	BMA	O5-C5-C6-O6
4	E	3	BMA	C4-C5-C6-O6
4	D	3	BMA	C4-C5-C6-O6
5	F	2	NAG	O5-C5-C6-O6
4	E	1	NAG	C4-C5-C6-O6
4	D	1	NAG	C8-C7-N2-C2
4	D	1	NAG	O7-C7-N2-C2
5	F	1	NAG	C8-C7-N2-C2
5	F	1	NAG	O7-C7-N2-C2
4	D	1	NAG	O5-C5-C6-O6
4	D	1	NAG	C4-C5-C6-O6
5	F	2	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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





5.6 Ligand geometry [i](#)

7 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	$\# Z > 2$
6	NAG	B	1302	1	14,14,15	0.33	0	17,19,21	0.64	0
6	NAG	A	1302	1	14,14,15	0.29	0	17,19,21	0.75	0
6	NAG	C	5003	1	14,14,15	0.28	0	17,19,21	0.79	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	C	5001	1	14,14,15	0.30	0	17,19,21	0.62	0
6	NAG	C	5002	1	14,14,15	0.27	0	17,19,21	0.86	1 (5%)
6	NAG	B	1301	1	14,14,15	0.28	0	17,19,21	0.67	0
6	NAG	A	1301	1	14,14,15	0.33	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
6	NAG	B	1302	1	-	0/6/23/26	0/1/1/1
6	NAG	A	1302	1	-	0/6/23/26	0/1/1/1
6	NAG	C	5003	1	-	2/6/23/26	0/1/1/1
6	NAG	C	5001	1	-	2/6/23/26	0/1/1/1
6	NAG	C	5002	1	-	2/6/23/26	0/1/1/1
6	NAG	B	1301	1	-	2/6/23/26	0/1/1/1
6	NAG	A	1301	1	-	3/6/23/26	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(°)
6	C	5002	NAG	O5-C1-C2	-2.42	107.55	111.29

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	5001	NAG	O5-C5-C6-O6
6	A	1301	NAG	O5-C5-C6-O6
6	C	5002	NAG	O5-C5-C6-O6
6	C	5001	NAG	C4-C5-C6-O6
6	B	1301	NAG	O5-C5-C6-O6
6	A	1301	NAG	C4-C5-C6-O6
6	B	1301	NAG	C4-C5-C6-O6
6	C	5002	NAG	C4-C5-C6-O6
6	C	5003	NAG	C4-C5-C6-O6
6	A	1301	NAG	C1-C2-N2-C7
6	C	5003	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	5002	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.

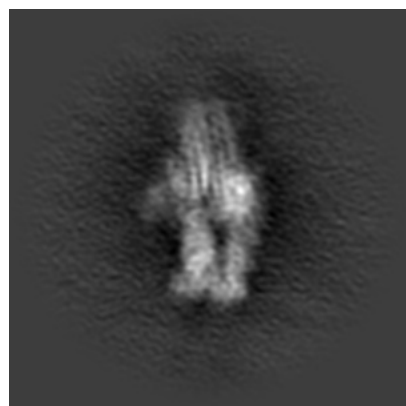
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23064. These allow visual inspection of the internal detail of the map and identification of artifacts.

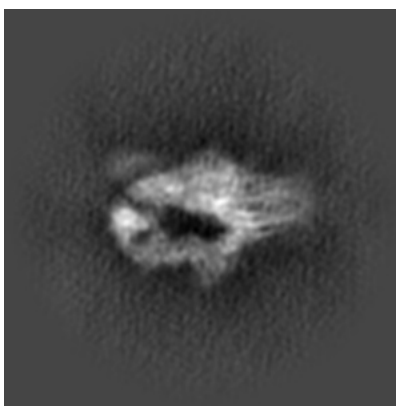
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

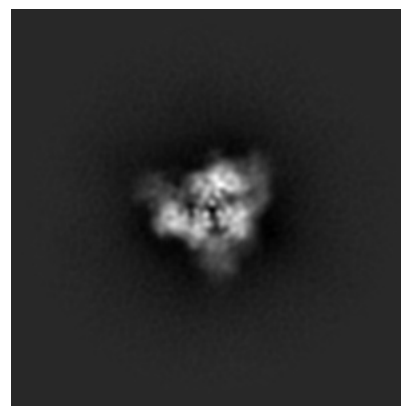
6.1.1 Primary map



X

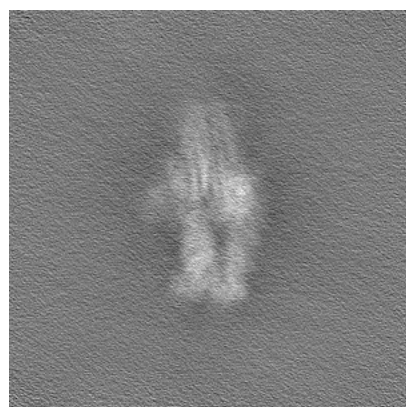


Y

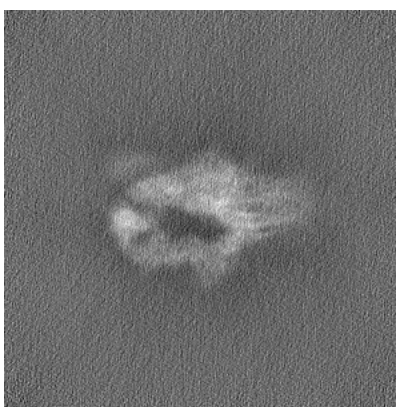


Z

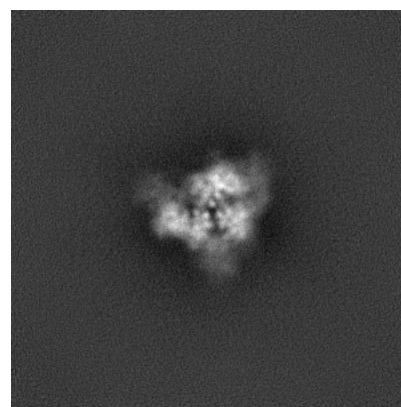
6.1.2 Raw map



X



Y

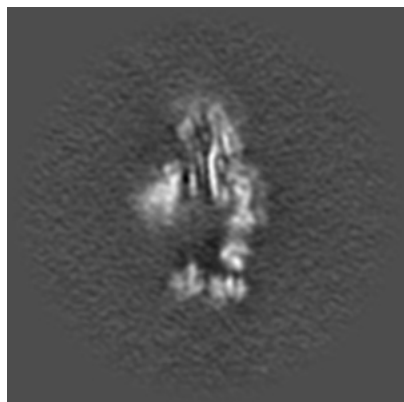


Z

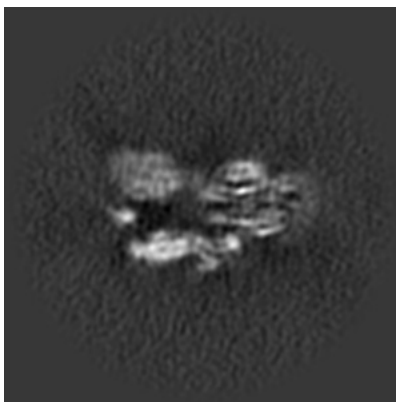
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

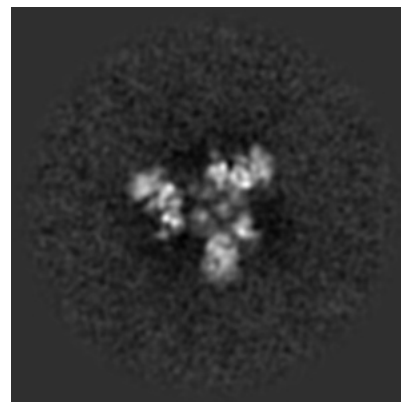
6.2.1 Primary map



X Index: 200

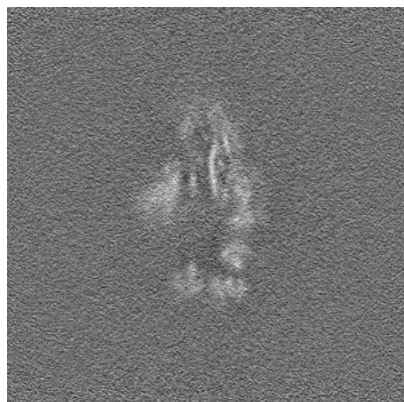


Y Index: 200



Z Index: 200

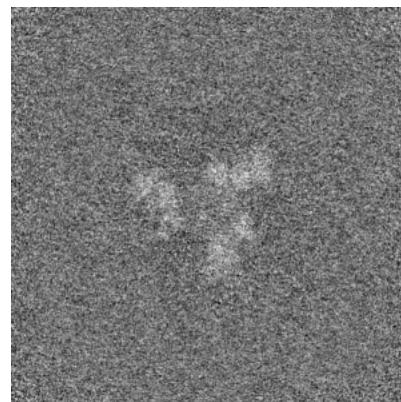
6.2.2 Raw map



X Index: 200



Y Index: 200

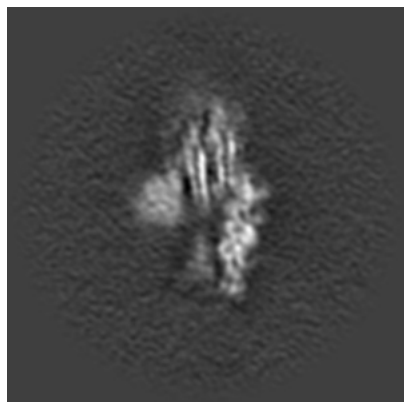


Z Index: 200

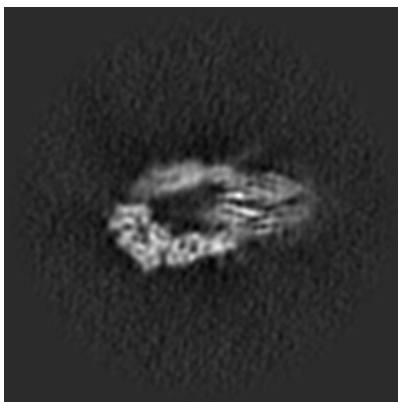
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

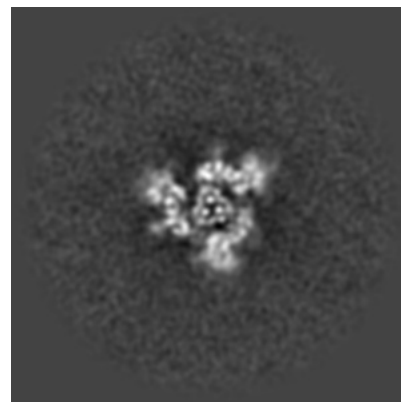
6.3.1 Primary map



X Index: 208

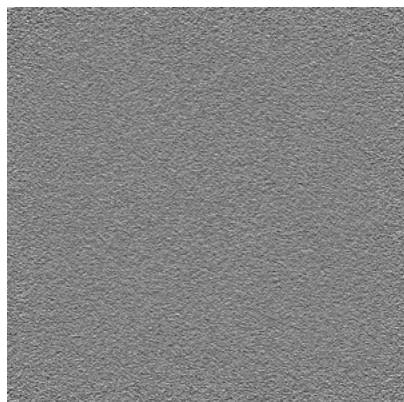


Y Index: 185

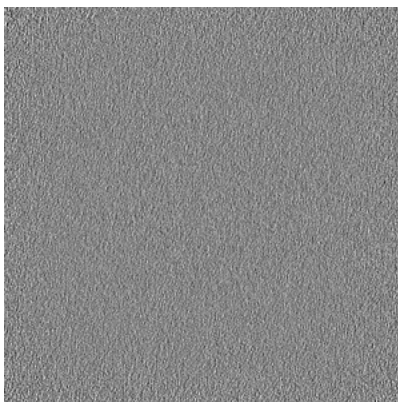


Z Index: 218

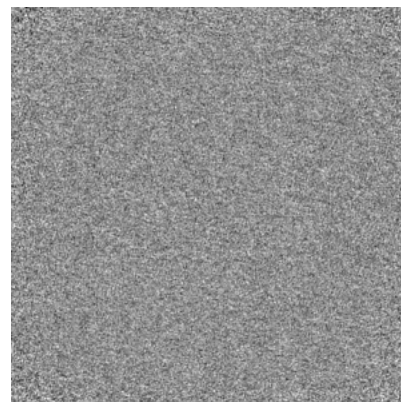
6.3.2 Raw map



X Index: 0



Y Index: 0

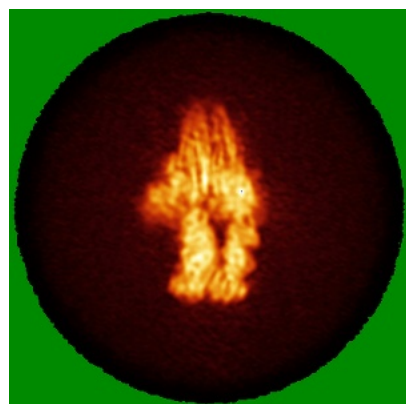


Z Index: 0

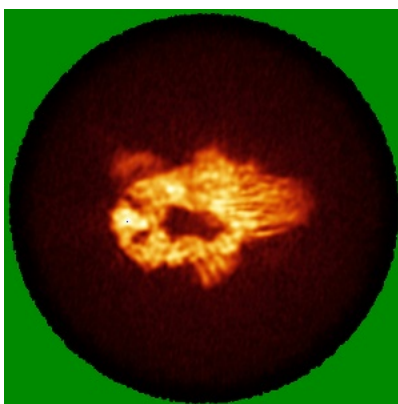
The images above show the largest variance slices of the map in three orthogonal directions.

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

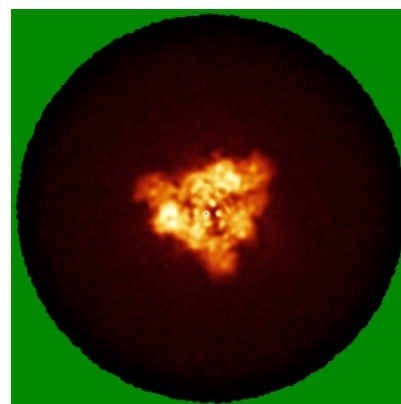
6.4.1 Primary map



X

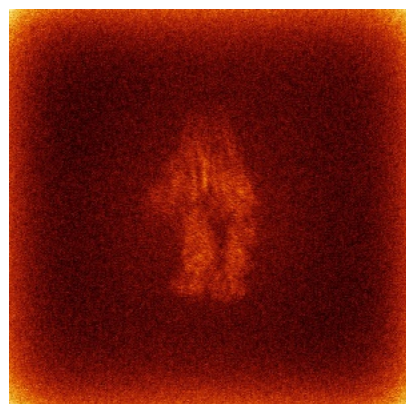


Y

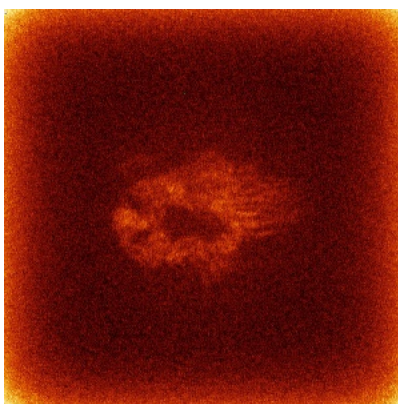


Z

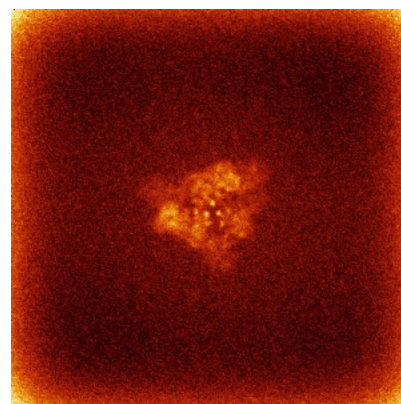
6.4.2 Raw map



X



Y

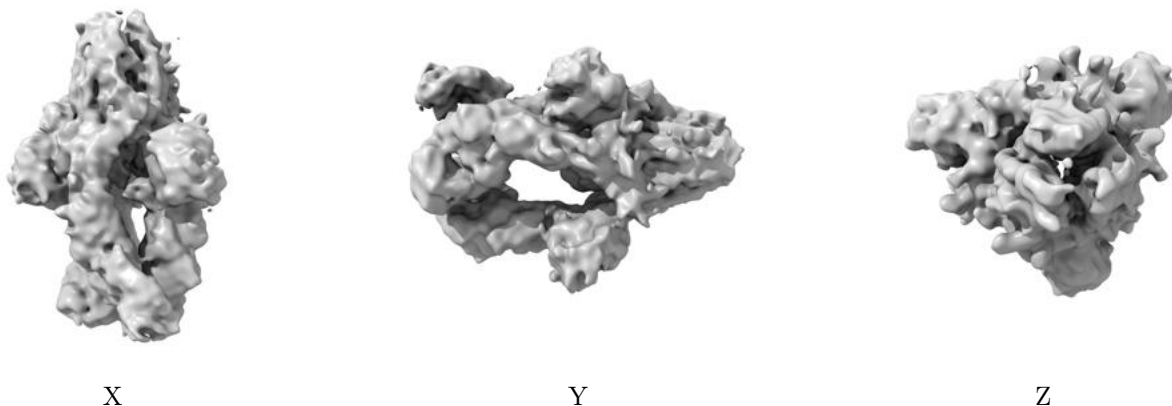


Z

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

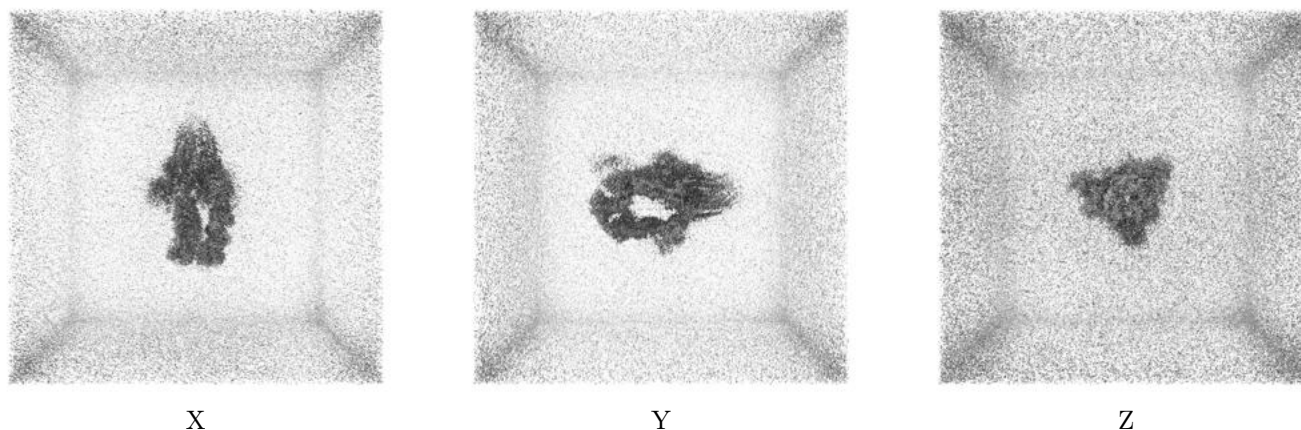
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.24. 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.

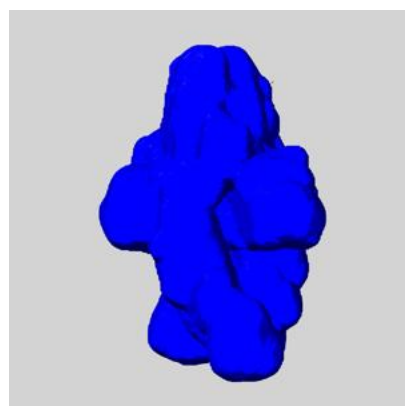
6.6 Mask visualisation [i](#)

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

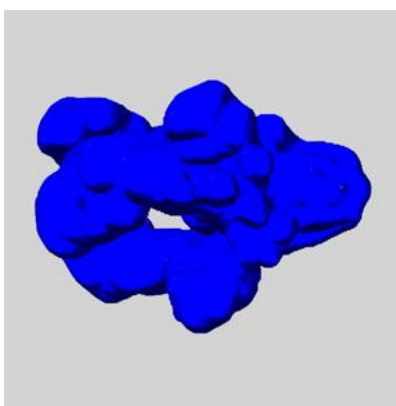
A mask typically either:

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

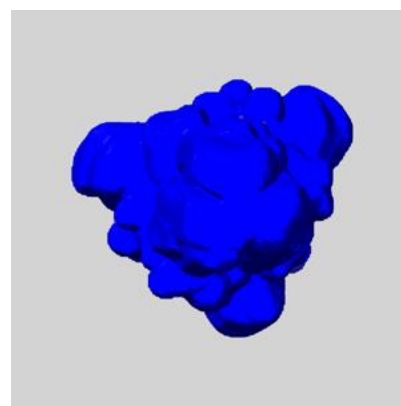
6.6.1 emd_23064_msk_1.map [i](#)



X



Y

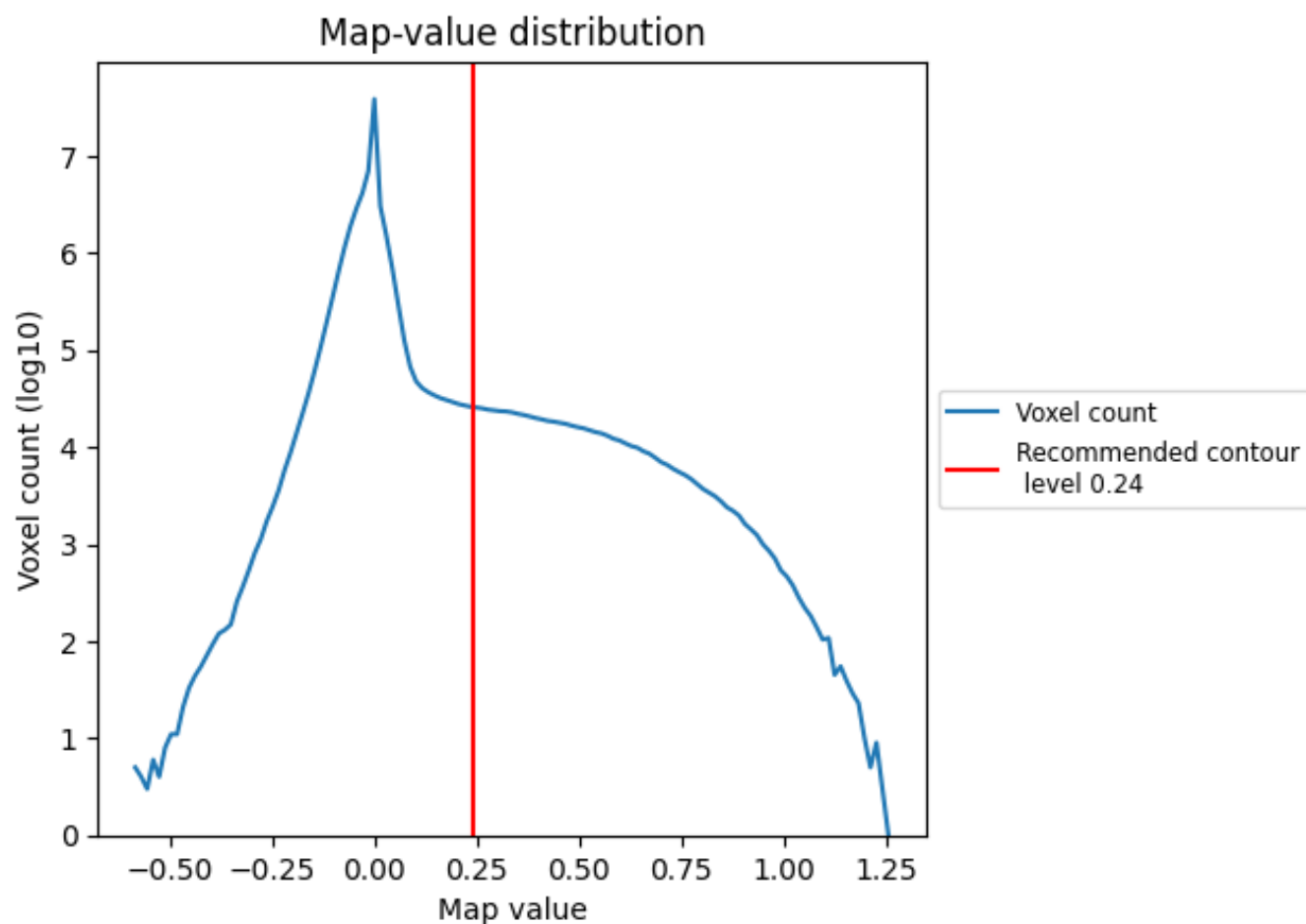


Z

7 Map analysis [i](#)

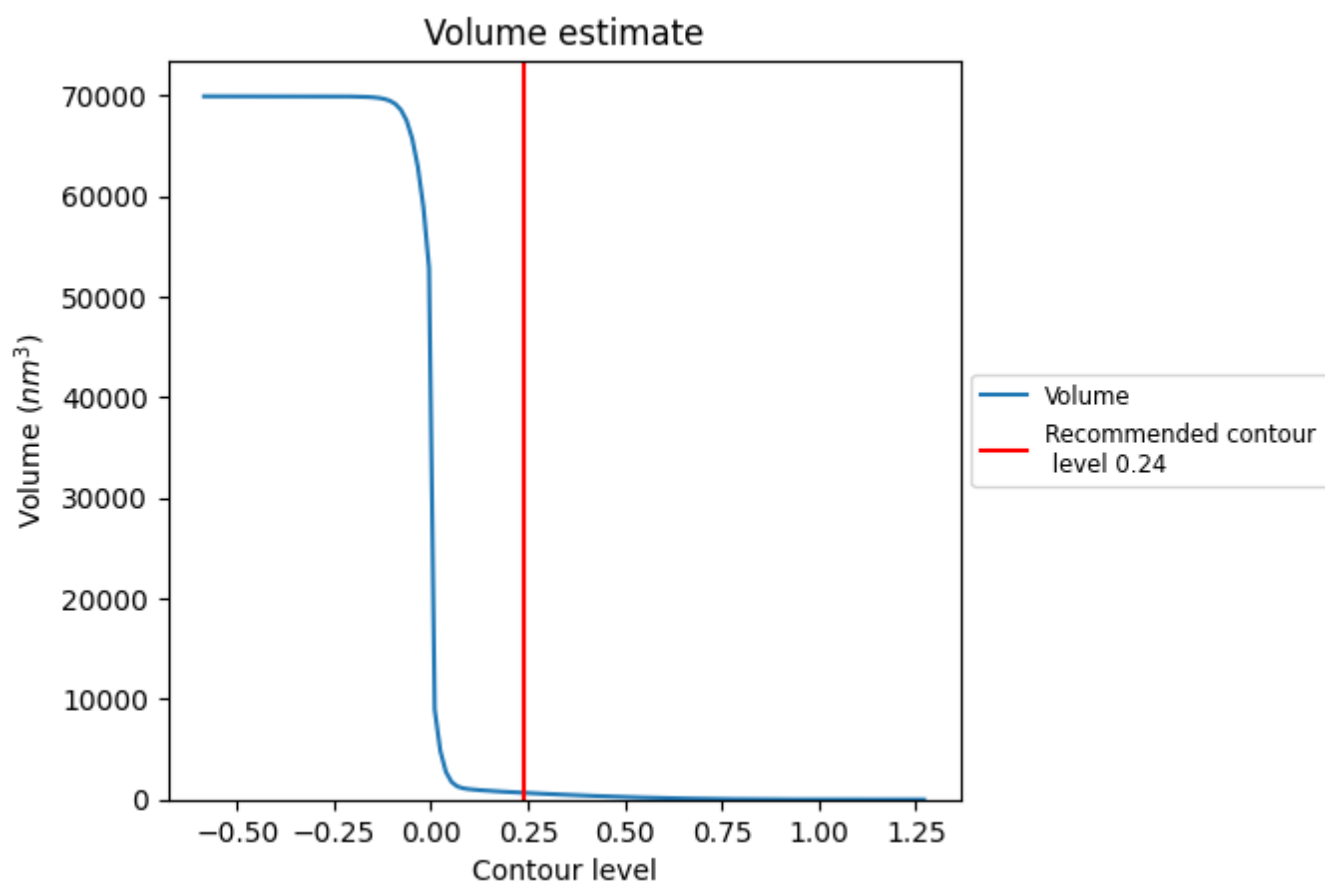
This section contains the results of statistical analysis of the map.

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.

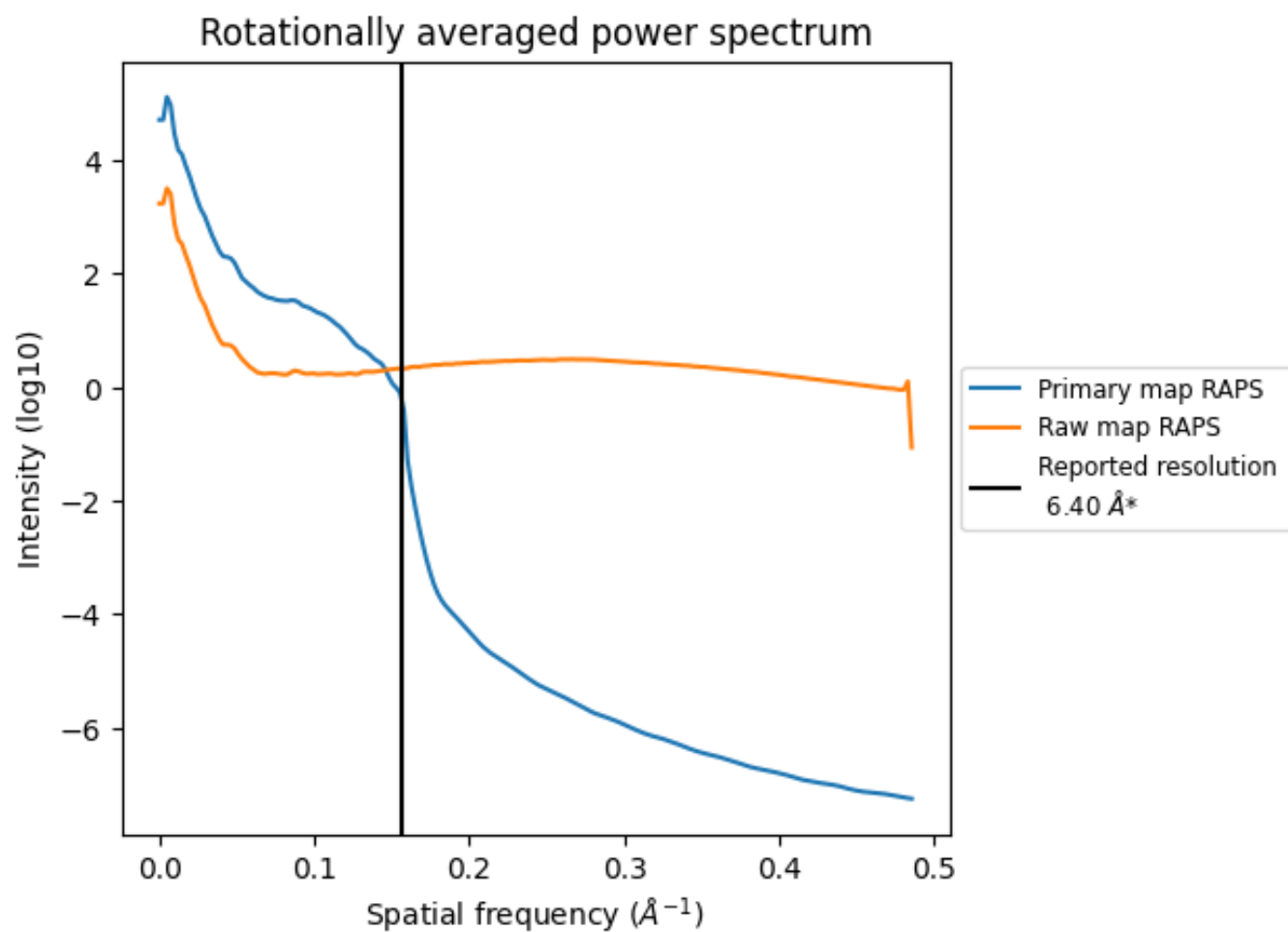
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 669 nm³; this corresponds to an approximate mass of 604 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 ⓘ

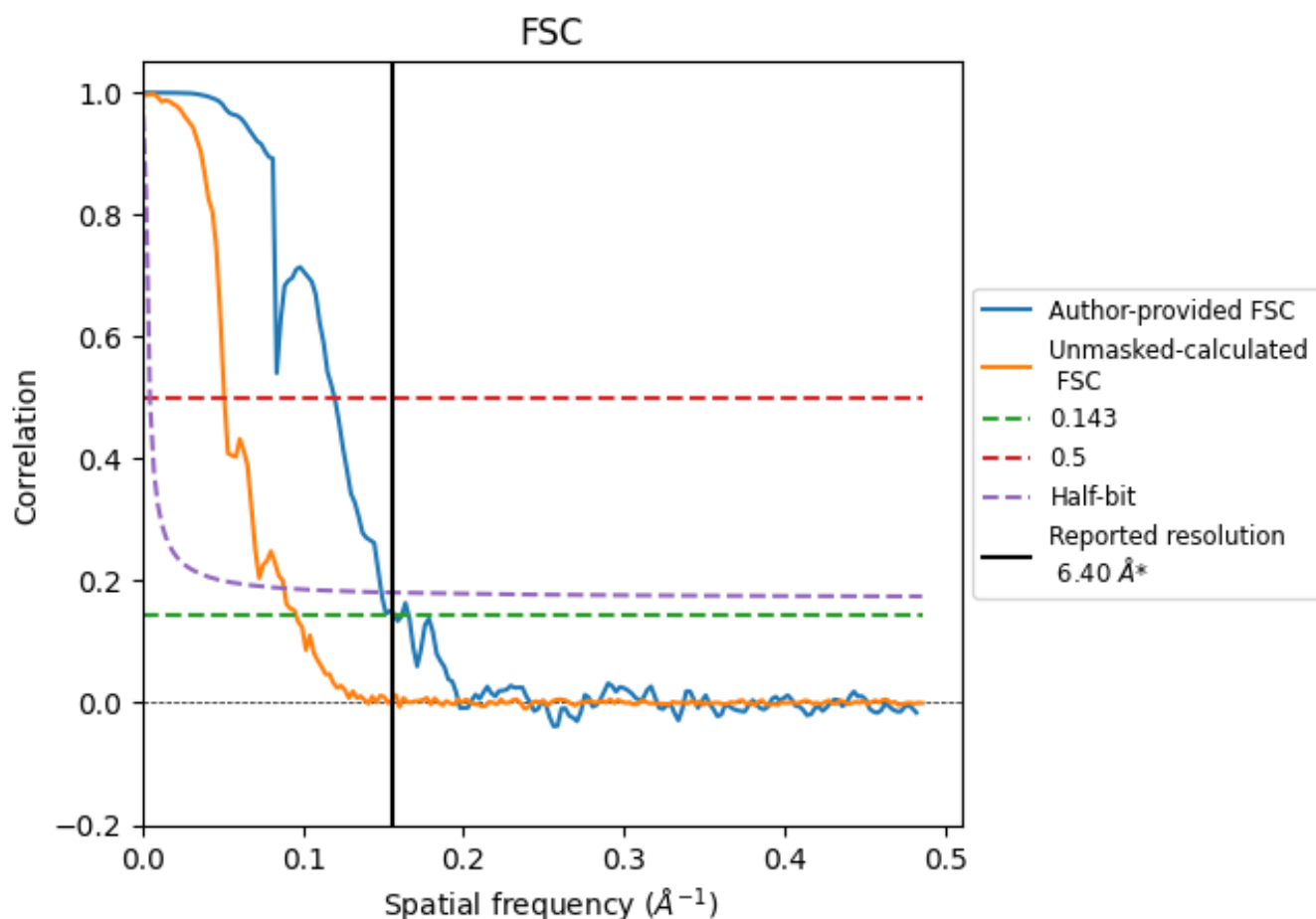


*Reported resolution corresponds to spatial frequency of 0.156 Å⁻¹

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 0.156 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

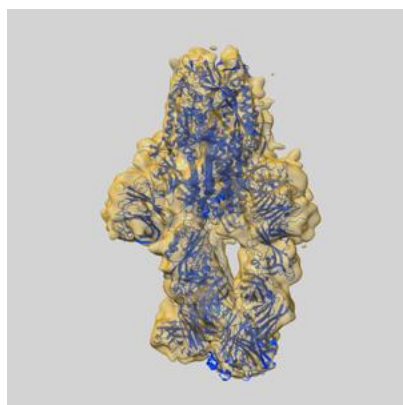
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	6.40	-	-
Author-provided FSC curve	6.36	8.35	6.72
Unmasked-calculated*	10.44	19.61	11.34

*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 10.44 differs from the reported value 6.4 by more than 10 %

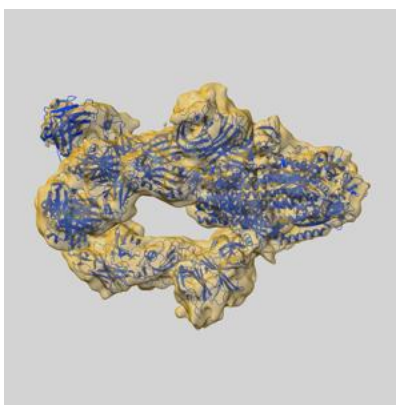
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23064 and PDB model 7KXJ. Per-residue inclusion information can be found in section 3 on page 12.

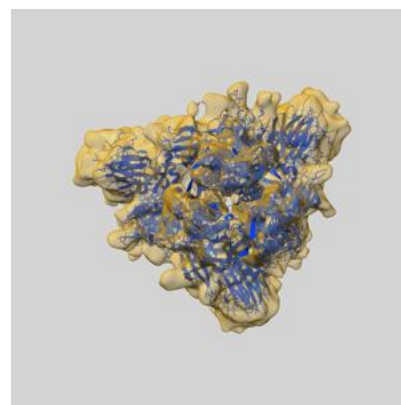
9.1 Map-model overlay [i](#)



X



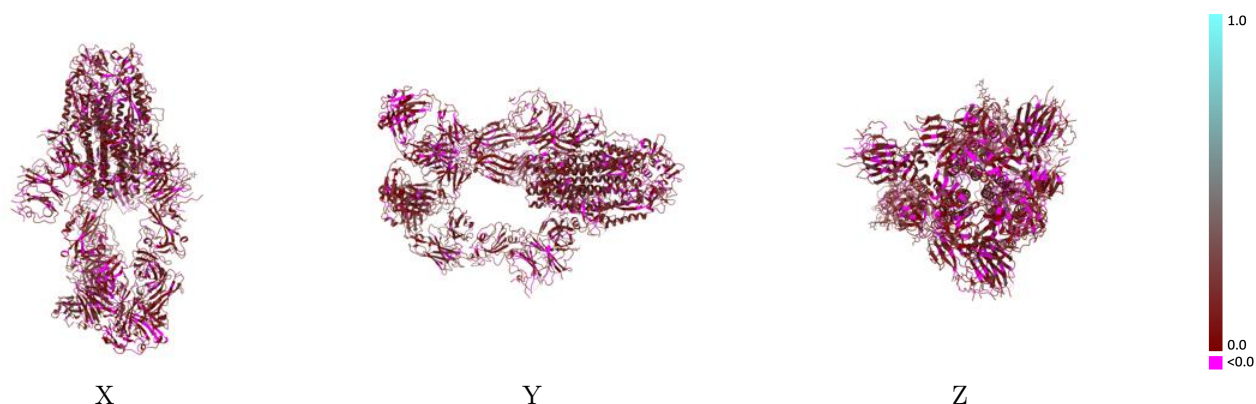
Y



Z

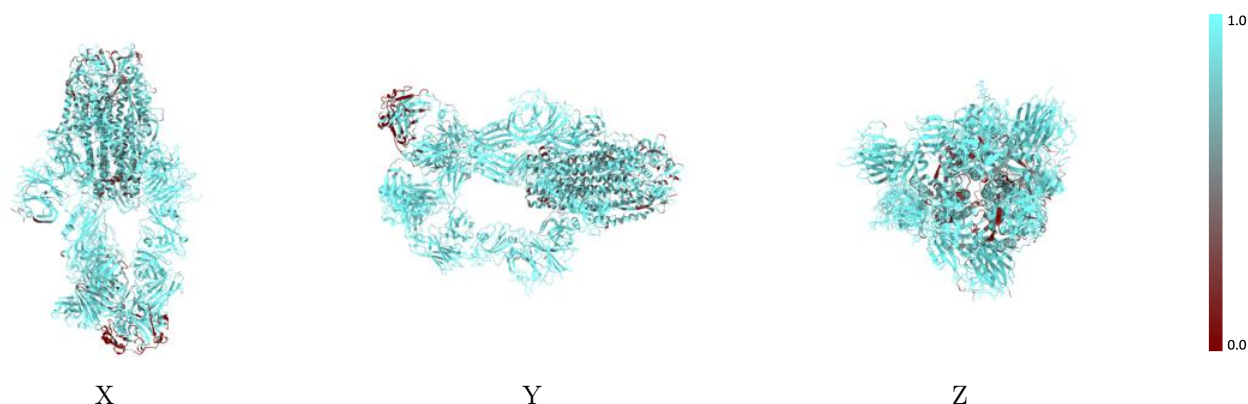
The images above show the 3D surface view of the map at the recommended contour level 0.24 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



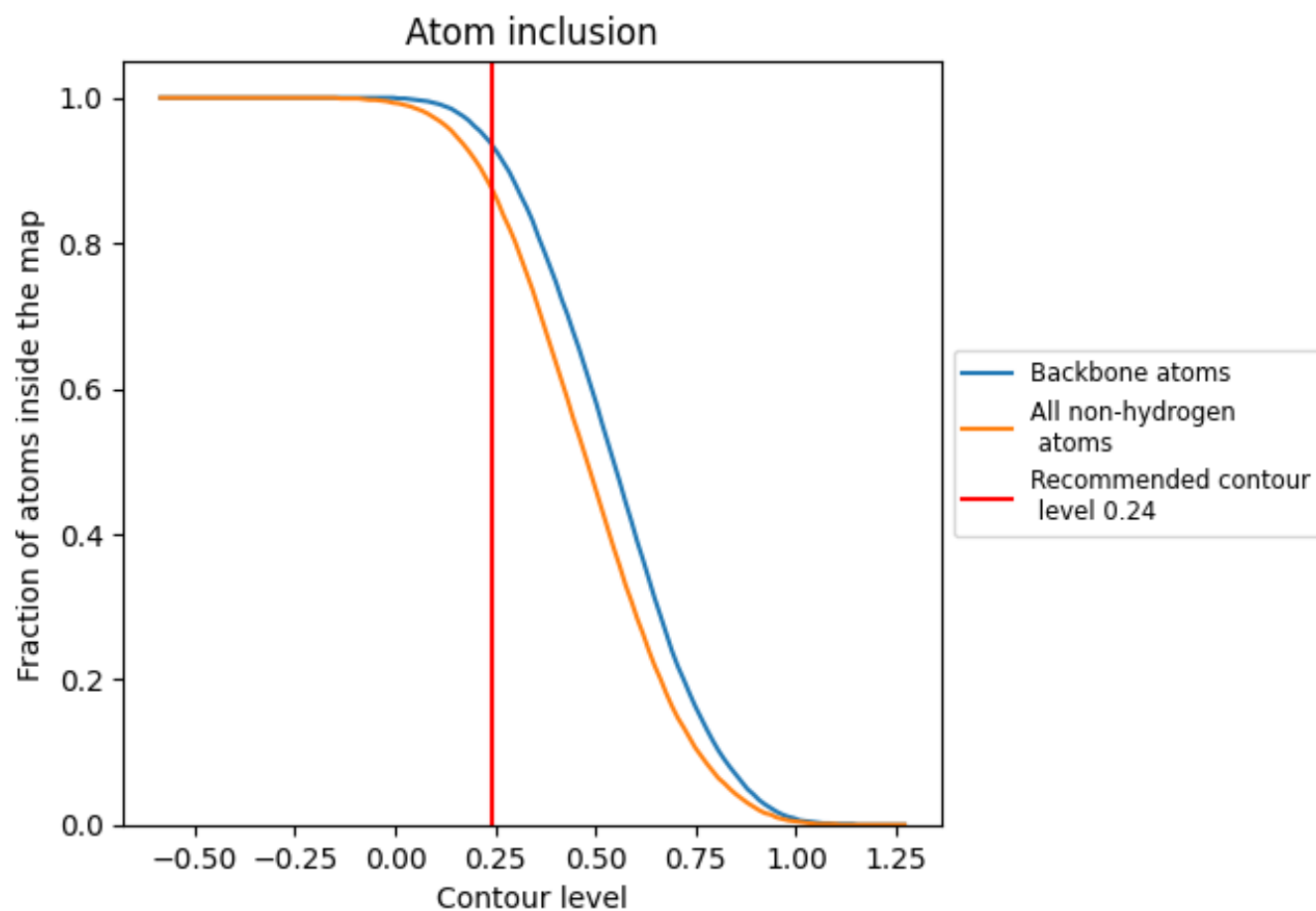
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.24).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 88% 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 (0.24) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8770	0.1130
A	0.8880	0.1150
B	0.8780	0.1140
C	0.8770	0.1090
D	0.8970	0.2470
E	0.9740	0.2830
F	1.0000	0.1440
H	0.9530	0.1250
I	0.9770	0.1410
J	0.5950	0.0720
L	0.9710	0.1280
M	0.9810	0.1440
N	0.7180	0.0650

