



Full wwPDB EM Validation Report ⓘ

Apr 5, 2026 – 04:10 PM UTC

PDB ID : 9FQT / pdb_00009fqt
EMDB ID : EMD-50668
Title : Cryo-EM structure of MmCAT1 bound with FrMLV-RBD in the apo inward-open state
Authors : Ye, M.; Zhou, D.; Pike, A.C.W.; Wang, S.; Wang, D.; Bakshi, S.; Brooke, L.; Williams, E.; Elkins, J.; Stuart, D.I.; Sauer, D.B.
Deposited on : 2024-06-17
Resolution : 3.50 Å (reported)
Based on initial models : ., 1AOL

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:

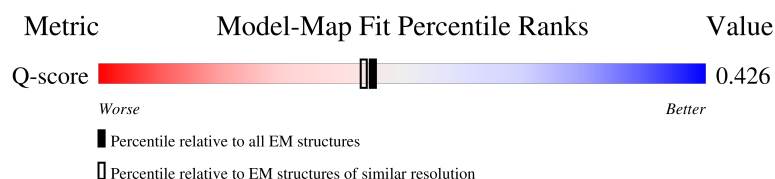
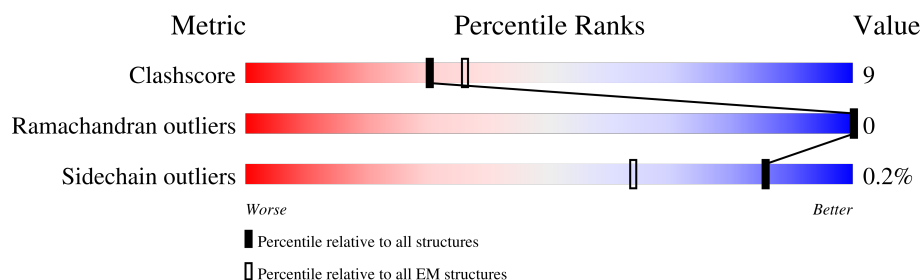
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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

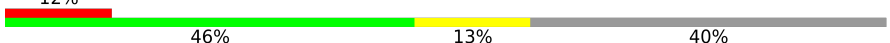
The reported resolution of this entry is 3.50 Å.

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	13950 (3.00 - 4.00)

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	902	
2	B	276	
3	D	5	

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Mol	Chain	Length	Quality of chain
4	C	2	50% 100%
4	I	2	50% 100%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11733 atoms, of which 5811 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 High affinity cationic amino acid transporter 1, Green fluorescent protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	537	Total	C	H	N	O	S	0	0
			8104	2654	4084	628	711	27		

There are 59 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	623	ALA	-	linker	UNP Q09143
A	624	GLU	-	linker	UNP Q09143
A	625	ASN	-	linker	UNP Q09143
A	626	LEU	-	linker	UNP Q09143
A	627	TYR	-	linker	UNP Q09143
A	628	PHE	-	linker	UNP Q09143
A	629	GLN	-	linker	UNP Q09143
A	630	SER	-	linker	UNP Q09143
A	631	GLY	-	linker	UNP Q09143
A	632	SER	-	linker	UNP Q09143
A	633	ALA	-	linker	UNP Q09143
A	634	VAL	-	linker	UNP Q09143
A	697	LEU	PHE	conflict	UNP P42212
A	698	THR	SER	conflict	UNP P42212
A	839	LYS	ALA	conflict	UNP P42212
A	864	LEU	HIS	conflict	UNP P42212
A	872	SER	-	expression tag	UNP P42212
A	873	GLY	-	expression tag	UNP P42212
A	874	LEU	-	expression tag	UNP P42212
A	875	ARG	-	expression tag	UNP P42212
A	876	SER	-	expression tag	UNP P42212
A	877	TRP	-	expression tag	UNP P42212
A	878	SER	-	expression tag	UNP P42212
A	879	HIS	-	expression tag	UNP P42212
A	880	PRO	-	expression tag	UNP P42212
A	881	GLN	-	expression tag	UNP P42212
A	882	PHE	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
A	883	GLU	-	expression tag	UNP P42212
A	884	LYS	-	expression tag	UNP P42212
A	885	GLY	-	expression tag	UNP P42212
A	886	GLY	-	expression tag	UNP P42212
A	887	GLY	-	expression tag	UNP P42212
A	888	SER	-	expression tag	UNP P42212
A	889	GLY	-	expression tag	UNP P42212
A	890	GLY	-	expression tag	UNP P42212
A	891	GLY	-	expression tag	UNP P42212
A	892	SER	-	expression tag	UNP P42212
A	893	GLY	-	expression tag	UNP P42212
A	894	GLY	-	expression tag	UNP P42212
A	895	SER	-	expression tag	UNP P42212
A	896	ALA	-	expression tag	UNP P42212
A	897	TRP	-	expression tag	UNP P42212
A	898	SER	-	expression tag	UNP P42212
A	899	HIS	-	expression tag	UNP P42212
A	900	PRO	-	expression tag	UNP P42212
A	901	GLN	-	expression tag	UNP P42212
A	902	PHE	-	expression tag	UNP P42212
A	903	GLU	-	expression tag	UNP P42212
A	904	LYS	-	expression tag	UNP P42212
A	905	HIS	-	expression tag	UNP P42212
A	906	HIS	-	expression tag	UNP P42212
A	907	HIS	-	expression tag	UNP P42212
A	908	HIS	-	expression tag	UNP P42212
A	909	HIS	-	expression tag	UNP P42212
A	910	HIS	-	expression tag	UNP P42212
A	911	HIS	-	expression tag	UNP P42212
A	912	HIS	-	expression tag	UNP P42212
A	913	HIS	-	expression tag	UNP P42212
A	914	HIS	-	expression tag	UNP P42212

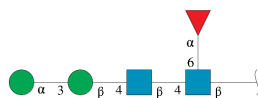
- Molecule 2 is a protein called Surface protein.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	B	223	Total	C	H	N	O	S	0	0
			3320	1095	1583	301	328	13		

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-30	MET	-	initiating methionine	UNP P03390
B	-29	GLY	-	expression tag	UNP P03390
B	-28	ILE	-	expression tag	UNP P03390
B	-27	LEU	-	expression tag	UNP P03390
B	-26	PRO	-	expression tag	UNP P03390
B	-25	SER	-	expression tag	UNP P03390
B	-24	PRO	-	expression tag	UNP P03390
B	-23	GLY	-	expression tag	UNP P03390
B	-22	MET	-	expression tag	UNP P03390
B	-21	PRO	-	expression tag	UNP P03390
B	-20	ALA	-	expression tag	UNP P03390
B	-19	LEU	-	expression tag	UNP P03390
B	-18	LEU	-	expression tag	UNP P03390
B	-17	SER	-	expression tag	UNP P03390
B	-16	LEU	-	expression tag	UNP P03390
B	-15	VAL	-	expression tag	UNP P03390
B	-14	SER	-	expression tag	UNP P03390
B	-13	LEU	-	expression tag	UNP P03390
B	-12	LEU	-	expression tag	UNP P03390
B	-11	SER	-	expression tag	UNP P03390
B	-10	VAL	-	expression tag	UNP P03390
B	-9	LEU	-	expression tag	UNP P03390
B	-8	LEU	-	expression tag	UNP P03390
B	-7	MET	-	expression tag	UNP P03390
B	-6	GLY	-	expression tag	UNP P03390
B	-5	CYS	-	expression tag	UNP P03390
B	-4	VAL	-	expression tag	UNP P03390
B	-3	ALA	-	expression tag	UNP P03390
B	-2	GLU	-	expression tag	UNP P03390
B	-1	THR	-	expression tag	UNP P03390
B	0	GLY	-	expression tag	UNP P03390
B	237	GLY	-	expression tag	UNP P03390
B	238	THR	-	expression tag	UNP P03390
B	239	LYS	-	expression tag	UNP P03390
B	240	HIS	-	expression tag	UNP P03390
B	241	HIS	-	expression tag	UNP P03390
B	242	HIS	-	expression tag	UNP P03390
B	243	HIS	-	expression tag	UNP P03390
B	244	HIS	-	expression tag	UNP P03390
B	245	HIS	-	expression tag	UNP P03390

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



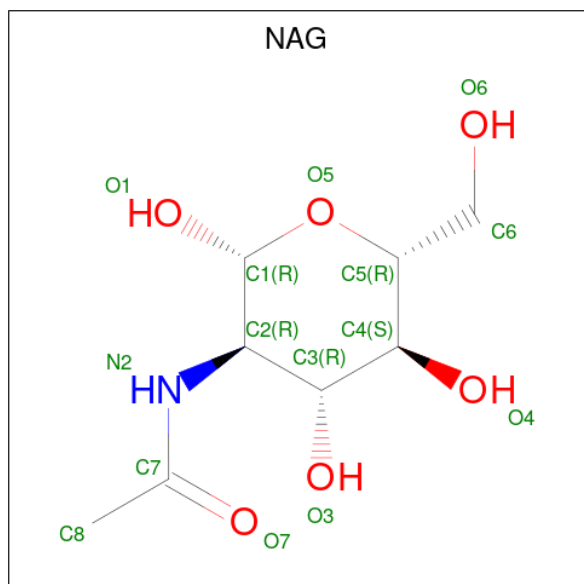
Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	5	Total	C	H	N	O	0	0
			97	34	37	2	24		

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



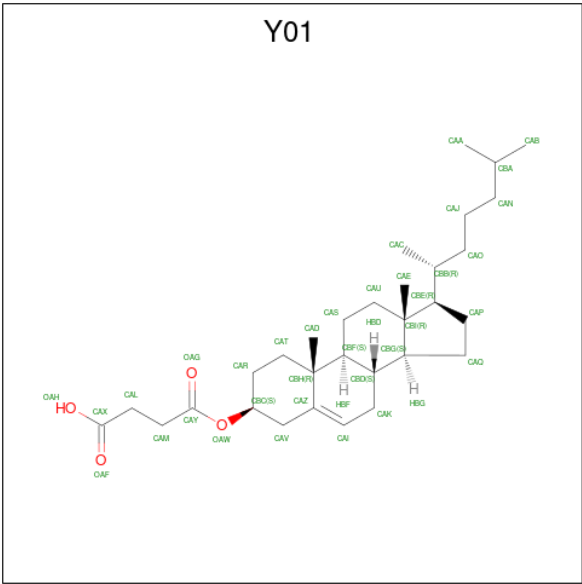
Mol	Chain	Residues	Atoms					AltConf	Trace
4	I	2	Total	C	H	N	O	0	0
			51	16	23	2	10		
4	C	2	Total	C	H	N	O	0	0
			51	16	23	2	10		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
5	A	1	Total	C	H	N	O	0
			26	8	12	1	5	

- Molecule 6 is CHOLESTEROL HEMISUCCINATE (CCD ID: Y01) (formula: $C_{31}H_{50}O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	H	O	
6	A	1	84	31	49	4	0

TYR	LEU	THR	GLN	SER	LYS	LEU	SER	ASP	LYS	PRO	ASN	GLU	LYS	ARG	ASP	HIS	MET	VAL	LEU	LEU	GLU	PHE	VAL	ALA	ALA	GLY	ILE	THR	LEU	GLY	MET	ASP	GLU	TYR	LYS	SER	GLY	LEU	ARG	SER	TRP	HIS	PRO	GLN	PHE	GLU	LYS	GLY	GLY	GLY	GLY	SER	GLY	GLY	GLY	SER																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																				
GLY	SER	ALA	TRP	HIS	PRO	GLN	PHE	LYS	GLU	LYS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS

• Molecule 2: Surface protein



PRO	GLY	THR	LYS	HIS	HIS	HIS	HIS	HIS	HIS																				Q9	V10	N19	E23	T24	V40	D44	L50	S51	P70	S76	S77	G78	S79			
L91	T95	N99	T100	M103	R104	L105	K106	L107	D108	H112	S115	E116	R128	E129	A130	K131	E147	V152	Y153	W154	D166	M167	T171	V175	C178	K179	D180	K196	S200	V213	S214	G215	R216	R225	L226	N230	L231	GLY	PRO	ARG	VAL				
MET	GLY	ILE	LEU	PRO	SER	PRO	GLY	MET	PRO	ALA	LEU	SER	VAL	LEU	LEU	MET	GLY	CYS	VAL	ALA	GLU	THR	GLY	ALA	ALA	PRO	GLY	SER	SER	PRO	HIS	Q9	V10	N19	E23	T24	V40	D44	L50	S51	P70	S76	S77	G78	S79

• Molecule 3: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1	MAG2	BMA3	MAN4	FUC5
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• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1	MAG2
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• Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1	MAG2
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	263131	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2600	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.251	Depositor
Minimum map value	-0.807	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.023	Depositor
Recommended contour level	0.197	Depositor
Map size (Å)	267.3, 267.3, 267.3	wwPDB
Map dimensions	324, 324, 324	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.825, 0.825, 0.825	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, FUC, Y01, 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	0.17	0/4114	0.36	0/5612
2	B	0.11	0/1799	0.31	0/2470
All	All	0.16	0/5913	0.34	0/8082

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	4020	4084	4096	90	0
2	B	1737	1583	1594	20	0
3	D	60	37	52	1	0
4	C	28	23	25	0	0
4	I	28	23	25	0	0
5	A	14	12	13	0	0
6	A	35	49	49	1	0
All	All	5922	5811	5854	106	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 9.

All (106) 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:496:LEU:O	1:A:500:ILE:HD12	1.73	0.89
3:D:2:NAG:O3	3:D:3:BMA:H2	1.80	0.82
2:B:166:ASP:OD1	2:B:167:ASN:N	2.14	0.79
1:A:117:LEU:HD22	1:A:354:VAL:HG21	1.72	0.71
1:A:360:GLU:OE2	1:A:369:ALA:HB1	1.91	0.71
1:A:52:VAL:HG22	1:A:398:LEU:HD21	1.75	0.69
2:B:147:GLU:N	2:B:147:GLU:OE1	2.25	0.69
1:A:216:ASN:ND2	1:A:303:LEU:O	2.27	0.67
1:A:75:LEU:HD13	6:A:1002:Y01:HAP1	1.75	0.67
1:A:52:VAL:CG2	1:A:398:LEU:HD21	2.26	0.66
1:A:41:GLY:HA3	1:A:265:ILE:HD11	1.79	0.64
1:A:176:THR:HG21	1:A:386:GLY:HA3	1.81	0.63
1:A:355:ILE:HG21	1:A:368:LEU:HD23	1.80	0.63
1:A:327:GLU:N	1:A:327:GLU:OE1	2.35	0.60
1:A:535:VAL:O	1:A:538:ILE:HG22	2.02	0.59
1:A:359:ALA:HB1	1:A:369:ALA:HB2	1.85	0.59
1:A:53:LEU:HD11	1:A:254:THR:HG22	1.86	0.58
1:A:201:LEU:HD22	1:A:297:VAL:HG21	1.86	0.56
1:A:52:VAL:HG21	1:A:131:ALA:HB2	1.88	0.55
1:A:48:ALA:HB1	1:A:340:LEU:HD23	1.89	0.55
1:A:237:GLU:OE2	2:B:95:THR:N	2.35	0.55
1:A:355:ILE:CG2	1:A:368:LEU:HD23	2.35	0.55
1:A:473:ILE:HG22	1:A:538:ILE:HD11	1.88	0.53
1:A:171:ILE:O	1:A:175:LEU:HD23	2.08	0.53
1:A:499:LEU:HD21	1:A:531:LEU:HD23	1.91	0.53
1:A:149:ARG:CZ	1:A:149:ARG:HA	2.39	0.53
1:A:124:GLY:O	1:A:128:VAL:HG23	2.09	0.53
2:B:50:LEU:O	2:B:51:SER:OG	2.24	0.53
1:A:235:TYR:CE2	2:B:70:PRO:HG3	2.44	0.52
1:A:50:VAL:HG13	1:A:51:TYR:HD2	1.75	0.51
2:B:100:THR:OG1	2:B:103:ASN:OD1	2.22	0.51
1:A:94:THR:HG21	1:A:269:GLY:HA2	1.93	0.51
1:A:144:ILE:H	1:A:144:ILE:HD12	1.76	0.51
1:A:584:VAL:O	1:A:588:ILE:HG13	2.10	0.51
1:A:221:GLU:OE1	2:B:99:ASN:ND2	2.41	0.50
1:A:103:VAL:HG23	1:A:104:THR:HG23	1.93	0.50
1:A:417:ALA:O	1:A:420:VAL:HG12	2.12	0.50
2:B:105:LEU:HD21	2:B:178:CYS:HB2	1.94	0.50
1:A:167:PHE:O	1:A:171:ILE:HG13	2.11	0.50
1:A:53:LEU:HD21	1:A:255:CYS:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:24:THR:OG1	2:B:225:ARG:NH2	2.45	0.49
1:A:304:MET:O	1:A:323:GLN:NE2	2.46	0.49
1:A:170:ILE:O	1:A:173:ILE:HG22	2.13	0.49
1:A:500:ILE:HG21	1:A:565:PHE:HA	1.95	0.49
1:A:425:TYR:CB	1:A:489:VAL:HG21	2.43	0.48
2:B:91:LEU:O	2:B:106:LYS:NZ	2.38	0.48
1:A:356:TYR:O	1:A:360:GLU:OE1	2.31	0.47
1:A:56:ALA:HB1	1:A:399:LYS:HE3	1.96	0.47
1:A:556:VAL:HG12	1:A:557:PRO:HD3	1.95	0.47
1:A:109:TRP:HZ2	1:A:415:LEU:O	1.97	0.47
1:A:101:SER:O	1:A:105:VAL:HG12	2.14	0.47
1:A:535:VAL:O	1:A:539:ILE:HG13	2.15	0.47
1:A:172:ILE:O	1:A:176:THR:HG23	2.15	0.46
1:A:404:LEU:O	1:A:407:ILE:HG22	2.15	0.46
1:A:485:SER:O	1:A:488:ILE:HG22	2.16	0.46
1:A:197:ASN:O	1:A:200:VAL:HG12	2.15	0.46
2:B:171:THR:O	2:B:175:VAL:HG23	2.15	0.46
1:A:142:LYS:O	1:A:146:GLU:N	2.44	0.46
1:A:425:TYR:HB3	1:A:489:VAL:HG21	1.97	0.46
1:A:48:ALA:HB3	1:A:343:SER:HB2	1.98	0.46
2:B:76:SER:OG	2:B:180:ASP:OD2	2.31	0.46
1:A:516:GLU:OE2	1:A:516:GLU:O	2.34	0.45
1:A:587:LEU:O	1:A:591:THR:HG23	2.16	0.45
1:A:428:GLU:HG2	1:A:428:GLU:O	2.16	0.45
1:A:117:LEU:HD22	1:A:354:VAL:CG2	2.44	0.44
1:A:500:ILE:HG23	1:A:565:PHE:HD1	1.81	0.44
1:A:33:ASN:C	1:A:33:ASN:HD22	2.26	0.44
1:A:499:LEU:O	1:A:503:VAL:HG23	2.18	0.44
1:A:267:THR:O	1:A:267:THR:HG22	2.17	0.44
1:A:359:ALA:CB	1:A:369:ALA:HB2	2.47	0.44
1:A:193:PHE:O	1:A:196:ILE:HG22	2.18	0.44
1:A:299:ALA:O	1:A:303:LEU:HG	2.17	0.43
2:B:40:VAL:HG23	2:B:40:VAL:O	2.17	0.43
1:A:149:ARG:HA	1:A:149:ARG:NE	2.34	0.43
2:B:213:VAL:HG12	2:B:214:SER:N	2.34	0.43
1:A:533:MET:SD	1:A:533:MET:C	3.02	0.43
2:B:108:ASP:OD1	2:B:112:HIS:CD2	2.71	0.43
1:A:356:TYR:CD1	1:A:360:GLU:OE1	2.72	0.43
1:A:94:THR:HG21	1:A:269:GLY:CA	2.48	0.42
1:A:480:GLU:HB2	1:A:481:PRO:HD2	2.01	0.42
1:A:90:ARG:NH2	1:A:549:LEU:HD21	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:TYR:OH	1:A:267:THR:HG23	2.20	0.42
1:A:516:GLU:OE1	2:B:115:SER:CB	2.68	0.42
1:A:105:VAL:HG13	1:A:109:TRP:HB3	2.01	0.42
2:B:19:ASN:ND2	2:B:51:SER:OG	2.52	0.42
1:A:119:LEU:HD23	1:A:119:LEU:C	2.45	0.42
1:A:260:VAL:HG13	1:A:413:TYR:OH	2.19	0.42
1:A:64:PRO:O	1:A:67:VAL:HG22	2.19	0.42
1:A:556:VAL:N	1:A:557:PRO:CD	2.83	0.42
1:A:51:TYR:OH	1:A:200:VAL:HG13	2.19	0.42
1:A:272:VAL:HG21	1:A:279:ILE:HD11	2.00	0.42
1:A:235:TYR:CD2	2:B:70:PRO:HG3	2.55	0.41
1:A:186:SER:HB2	1:A:189:VAL:HG12	2.02	0.41
1:A:512:GLU:OE2	1:A:512:GLU:N	2.47	0.41
1:A:162:GLN:N	1:A:397:GLU:OE2	2.53	0.41
1:A:233:VAL:HG21	1:A:235:TYR:CZ	2.55	0.41
1:A:121:TYR:HB3	1:A:351:MET:HB2	2.02	0.41
1:A:109:TRP:O	1:A:113:THR:HG23	2.21	0.41
1:A:193:PHE:HA	1:A:196:ILE:HG22	2.03	0.41
1:A:485:SER:O	1:A:489:VAL:HG23	2.20	0.41
1:A:57:VAL:HB	1:A:66:ILE:HD12	2.04	0.40
1:A:78:VAL:HG22	1:A:287:LEU:HD13	2.03	0.40
1:A:143:PRO:HA	1:A:146:GLU:HB2	2.03	0.40
1:A:275:PRO:O	1:A:279:ILE:HG12	2.21	0.40
2:B:226:LEU:HD23	2:B:226:LEU:C	2.46	0.40
2:B:152:VAL:HG21	2:B:154:TRP:CE2	2.57	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	533/902 (59%)	528 (99%)	5 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	221/276 (80%)	217 (98%)	4 (2%)	0	100	100
All	All	754/1178 (64%)	745 (99%)	9 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	423/749 (56%)	422 (100%)	1 (0%)	87	85
2	B	190/235 (81%)	190 (100%)	0	100	100
All	All	613/984 (62%)	612 (100%)	1 (0%)	85	85

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	257	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	155	ASN
1	A	232	ASN
1	A	426	GLN
1	A	573	GLN
2	B	19	ASN
2	B	61	GLN
2	B	112	HIS

5.3.3 RNA [i](#)

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 ⓘ

9 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	C	1	4,1	14,14,15	0.19	0	17,19,21	0.44	0
4	NAG	C	2	4	14,14,15	0.22	0	17,19,21	0.43	0
3	NAG	D	1	2,3	14,14,15	0.22	0	17,19,21	0.54	0
3	NAG	D	2	3	14,14,15	0.32	0	17,19,21	0.72	0
3	BMA	D	3	3	11,11,12	0.77	0	15,15,17	0.88	1 (6%)
3	MAN	D	4	3	11,11,12	0.54	0	15,15,17	0.97	2 (13%)
3	FUC	D	5	3	10,10,11	0.69	0	14,14,16	0.73	0
4	NAG	I	1	4,2	14,14,15	0.18	0	17,19,21	0.44	0
4	NAG	I	2	4	14,14,15	0.24	0	17,19,21	0.45	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	C	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	C	2	4	-	2/6/23/26	0/1/1/1
3	NAG	D	1	2,3	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	2/6/23/26	0/1/1/1
3	BMA	D	3	3	-	0/2/19/22	0/1/1/1
3	MAN	D	4	3	-	0/2/19/22	0/1/1/1
3	FUC	D	5	3	-	-	0/1/1/1
4	NAG	I	1	4,2	-	4/6/23/26	0/1/1/1
4	NAG	I	2	4	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	MAN	C1-O5-C5	2.53	115.58	112.19
3	D	4	MAN	O2-C2-C3	-2.19	105.61	110.15
3	D	3	BMA	O2-C2-C3	-2.02	105.96	110.15

There are no chirality outliers.

All (14) torsion outliers are listed below:

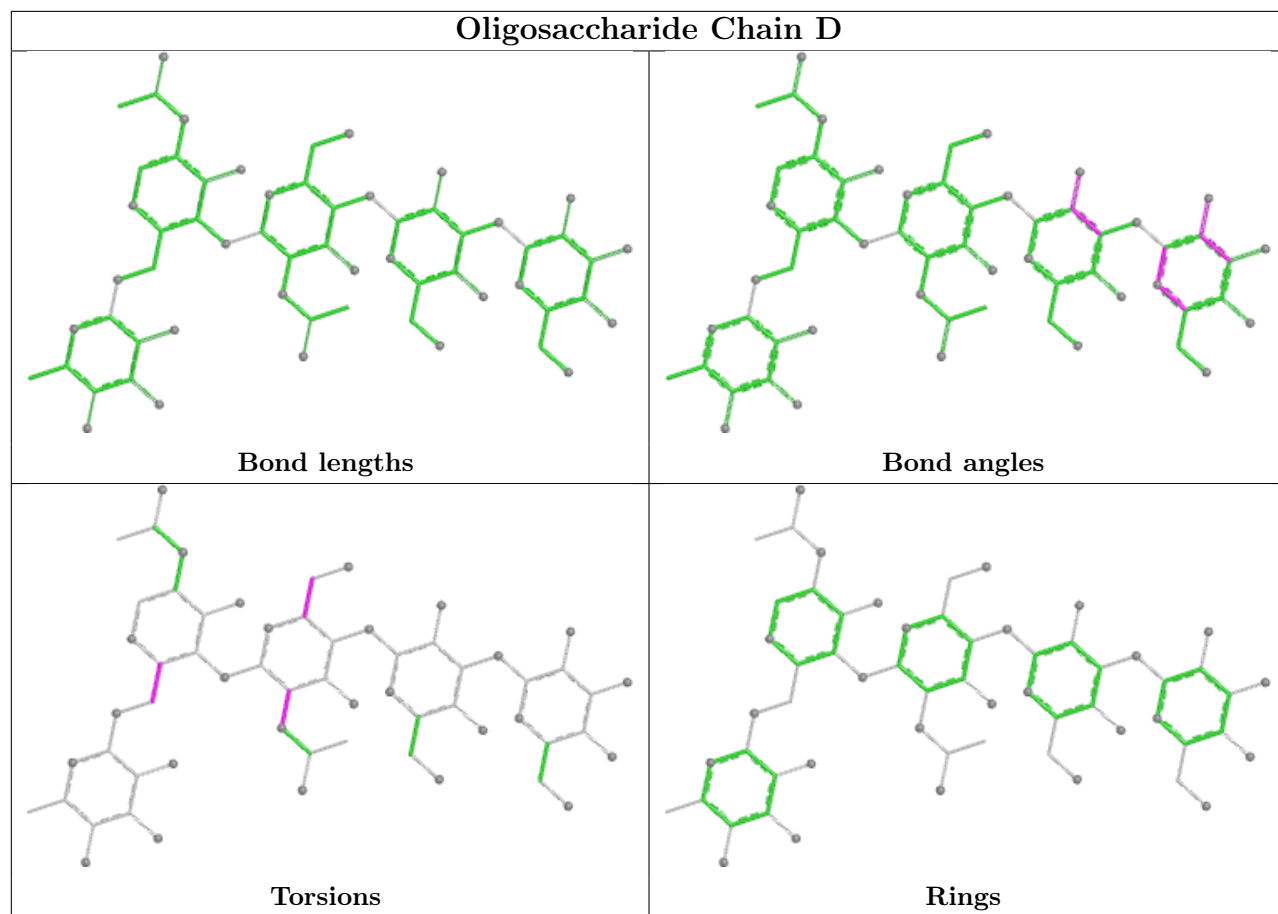
Mol	Chain	Res	Type	Atoms
4	C	2	NAG	O5-C5-C6-O6
4	C	2	NAG	C4-C5-C6-O6
4	C	1	NAG	O5-C5-C6-O6
4	C	1	NAG	C4-C5-C6-O6
3	D	2	NAG	C3-C2-N2-C7
3	D	1	NAG	C4-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
4	I	1	NAG	C1-C2-N2-C7
4	I	1	NAG	C3-C2-N2-C7
3	D	2	NAG	O5-C5-C6-O6
4	I	2	NAG	C4-C5-C6-O6
4	I	1	NAG	C4-C5-C6-O6
4	I	2	NAG	O5-C5-C6-O6
4	I	1	NAG	O5-C5-C6-O6

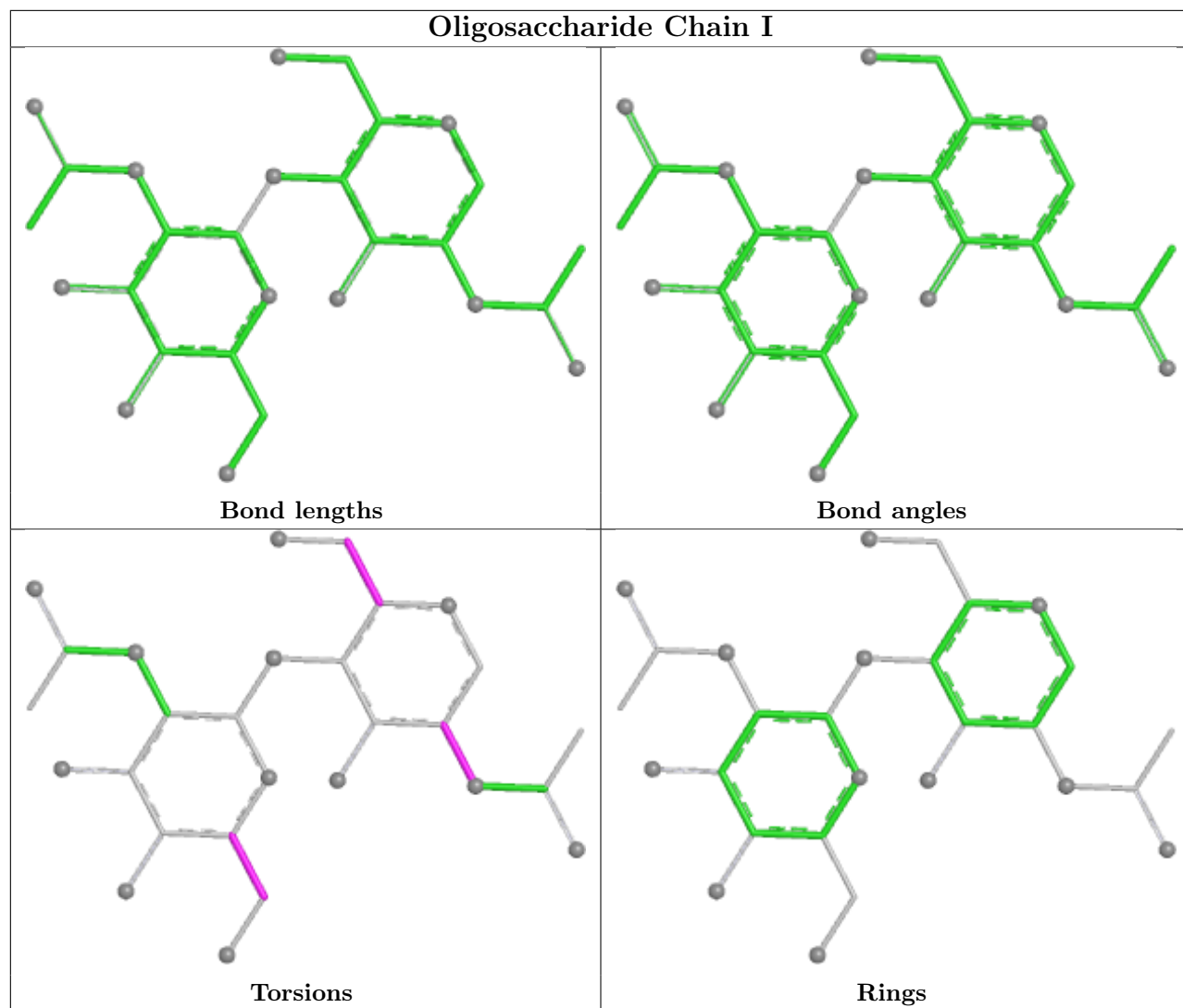
There are no ring outliers.

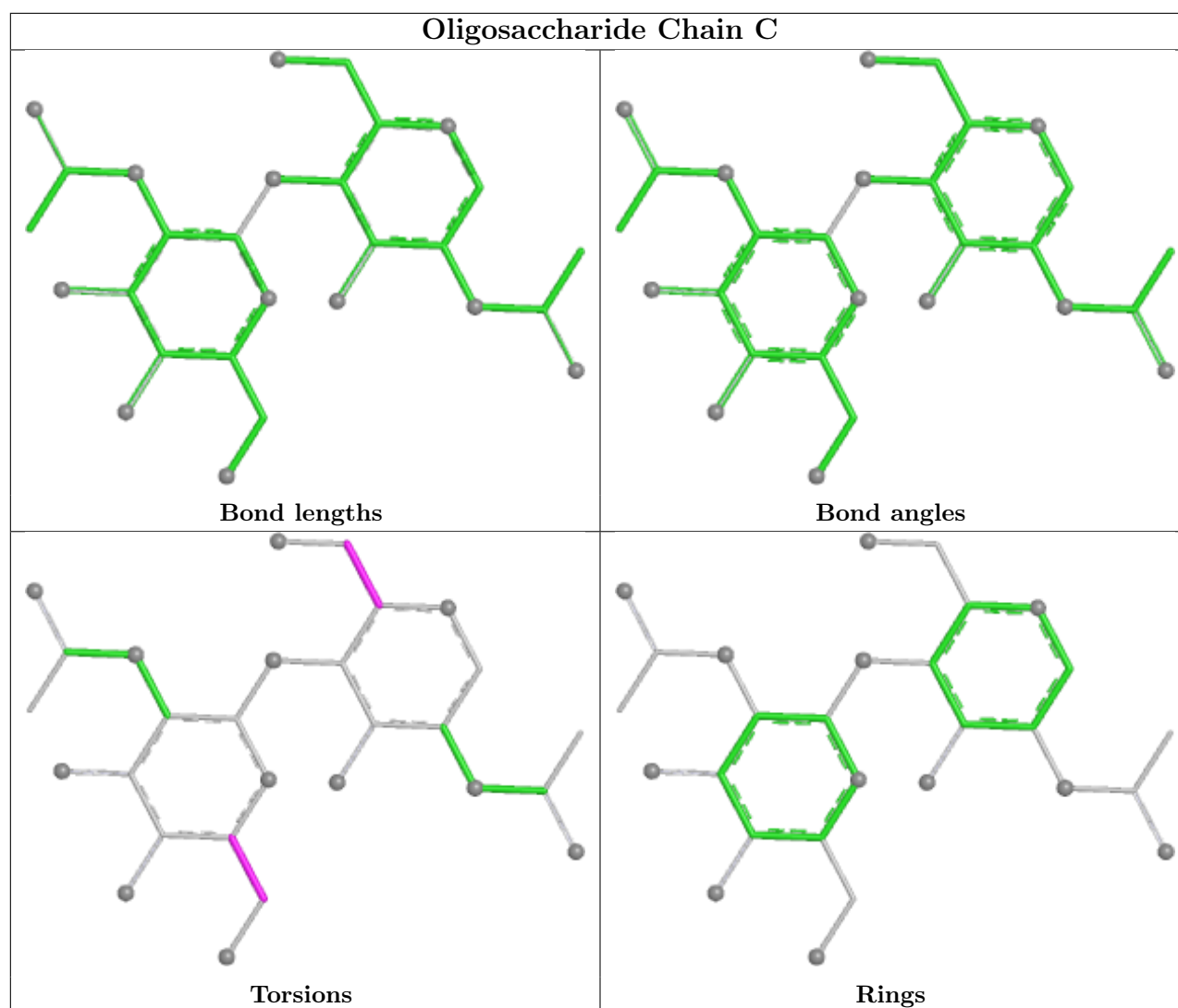
2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	2	NAG	1	0
3	D	3	BMA	1	0

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







5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
5	NAG	A	1001	1	14,14,15	0.32	0	17,19,21	0.56	0
6	Y01	A	1002	-	38,38,38	0.55	0	57,57,57	0.69	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1001	1	-	2/6/23/26	0/1/1/1
6	Y01	A	1002	-	-	2/19/77/77	0/4/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

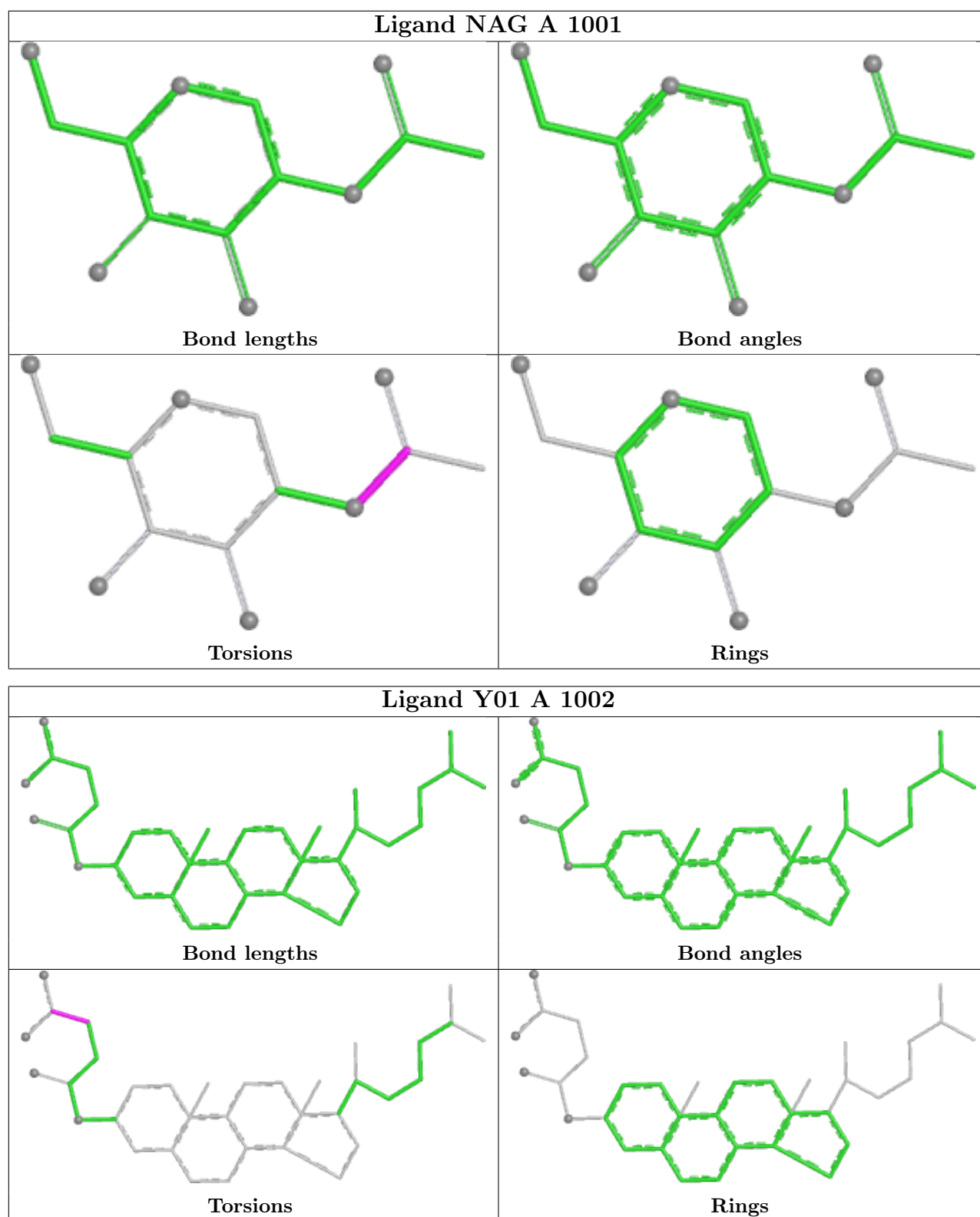
Mol	Chain	Res	Type	Atoms
5	A	1001	NAG	C8-C7-N2-C2
5	A	1001	NAG	O7-C7-N2-C2
6	A	1002	Y01	CAM-CAL-CAX-OAH
6	A	1002	Y01	CAM-CAL-CAX-OAF

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1002	Y01	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

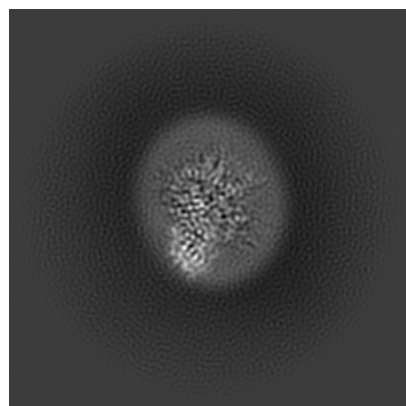
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-50668. These allow visual inspection of the internal detail of the map and identification of artifacts.

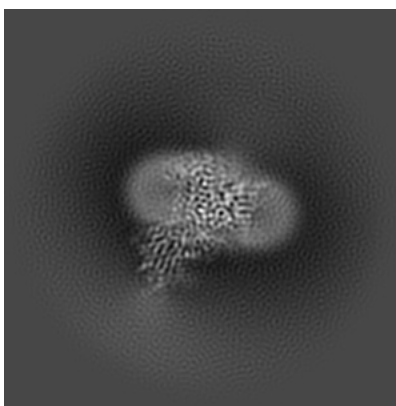
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

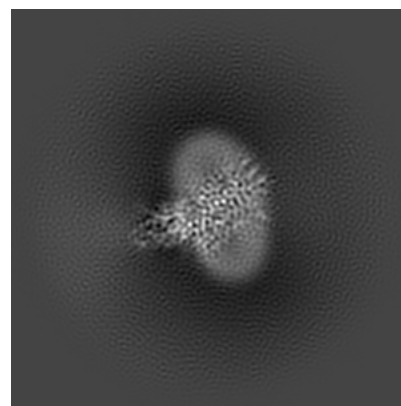
6.1.1 Primary map



X

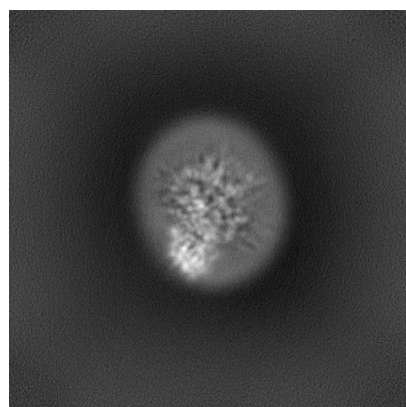


Y

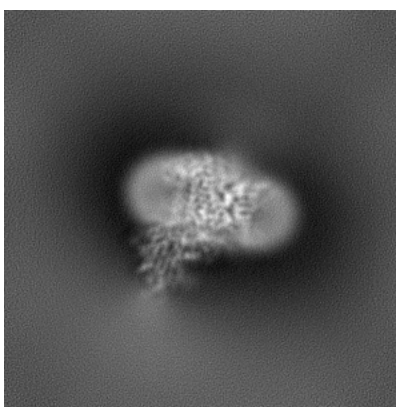


Z

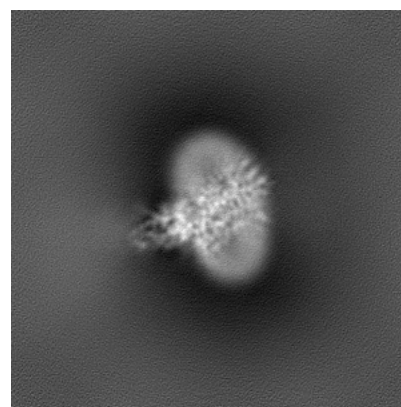
6.1.2 Raw map



X



Y

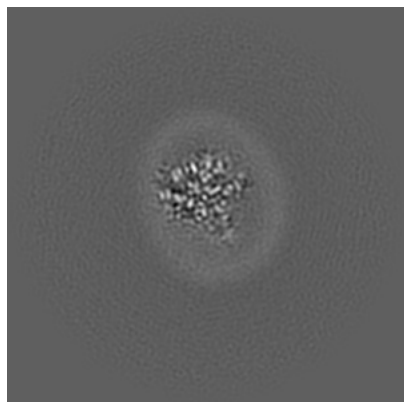


Z

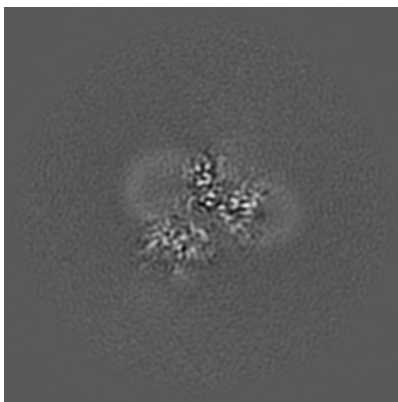
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

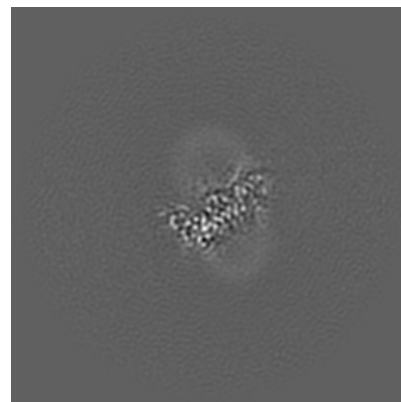
6.2.1 Primary map



X Index: 162

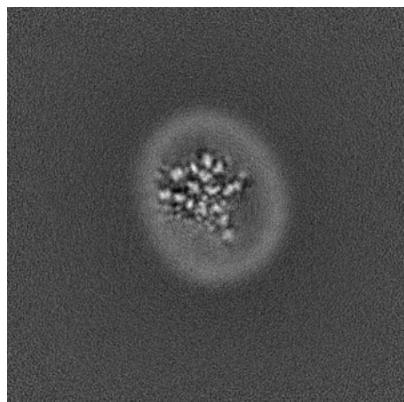


Y Index: 162

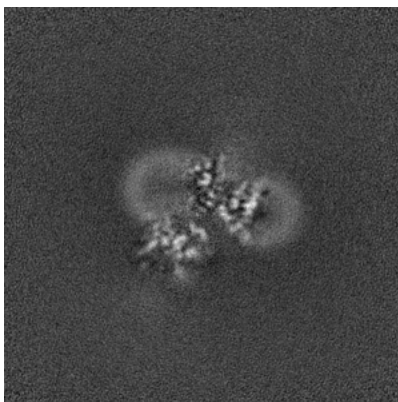


Z Index: 162

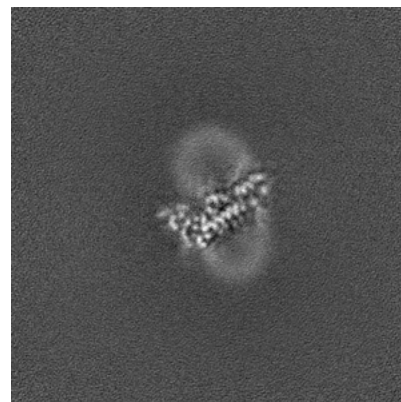
6.2.2 Raw map



X Index: 162



Y Index: 162

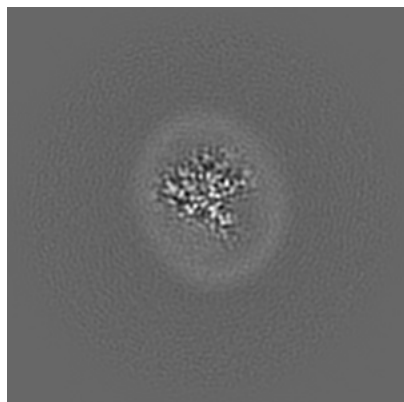


Z Index: 162

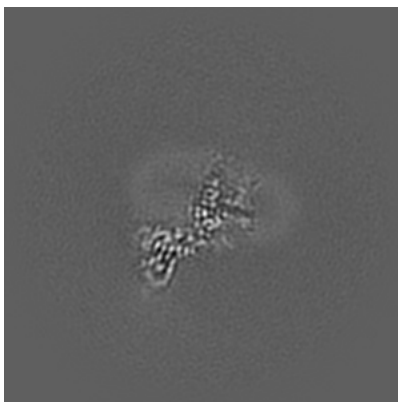
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

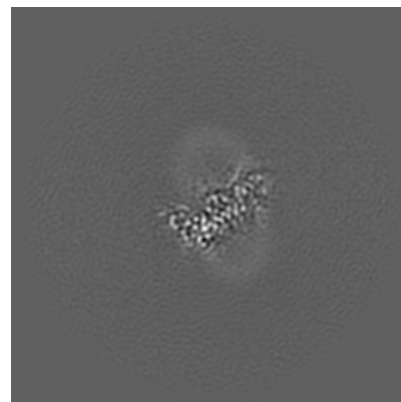
6.3.1 Primary map



X Index: 166

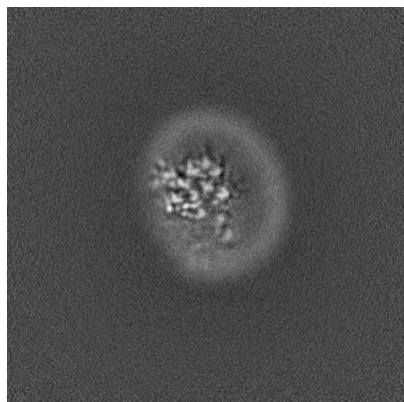


Y Index: 144

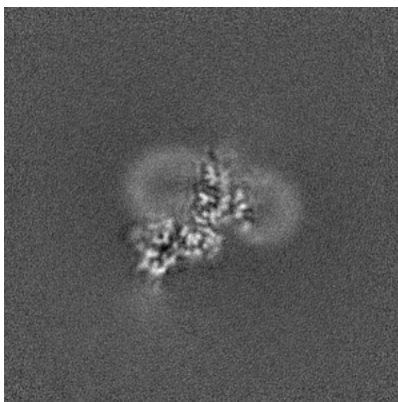


Z Index: 162

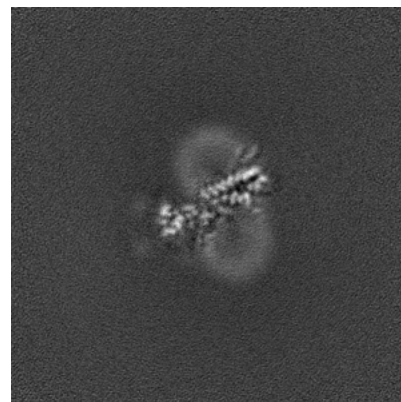
6.3.2 Raw map



X Index: 156



Y Index: 151

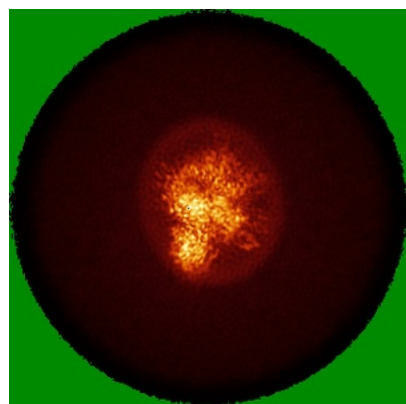


Z Index: 153

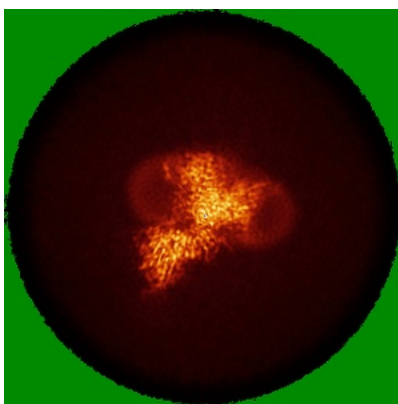
The images above show the largest variance slices of the map in three orthogonal directions.

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

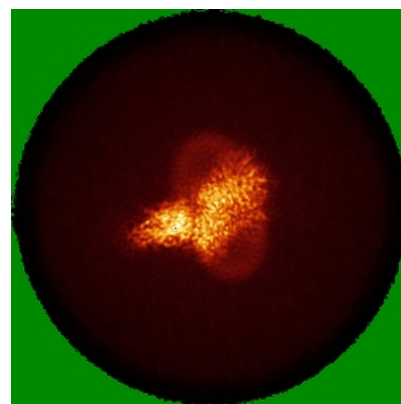
6.4.1 Primary map



X

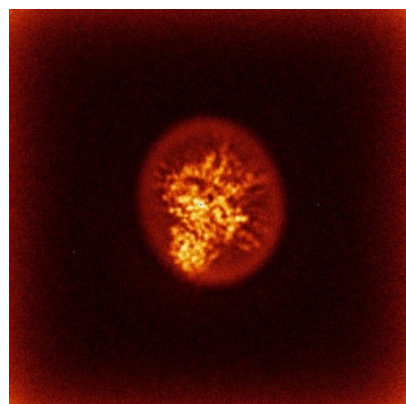


Y

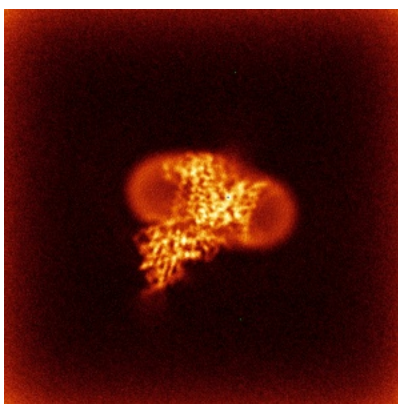


Z

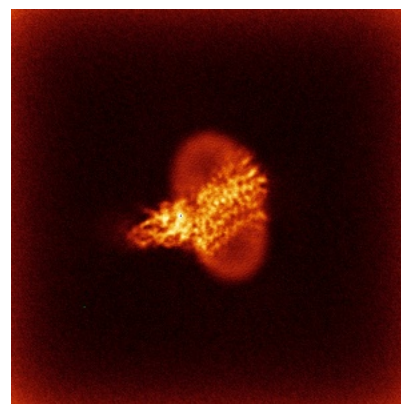
6.4.2 Raw map



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 0.197. 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.

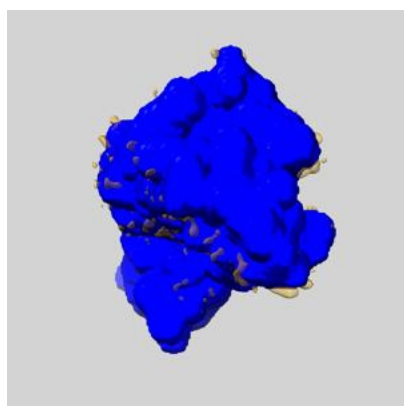
6.6 Mask visualisation [i](#)

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

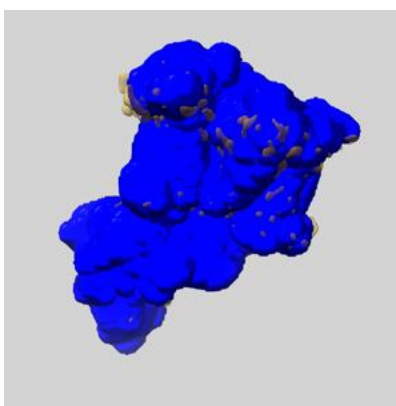
A mask typically either:

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

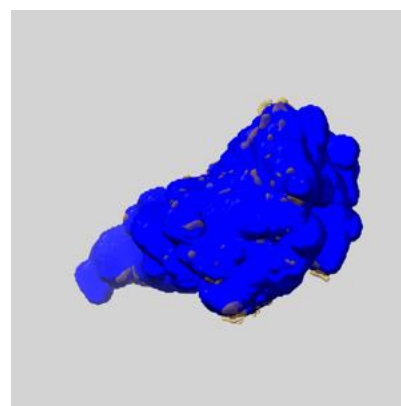
6.6.1 emd_50668_msk_1.map [i](#)



X



Y

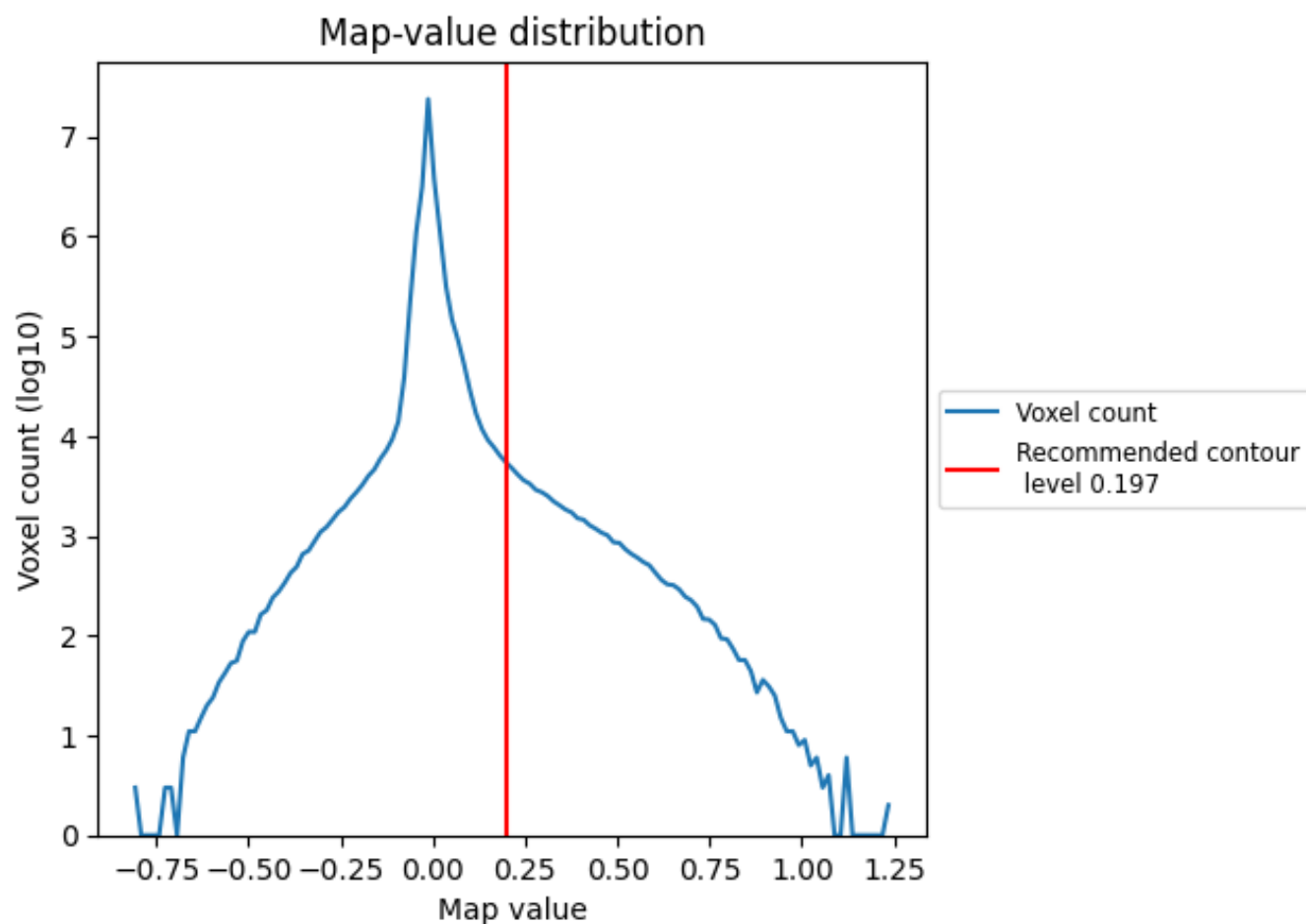


Z

7 Map analysis [i](#)

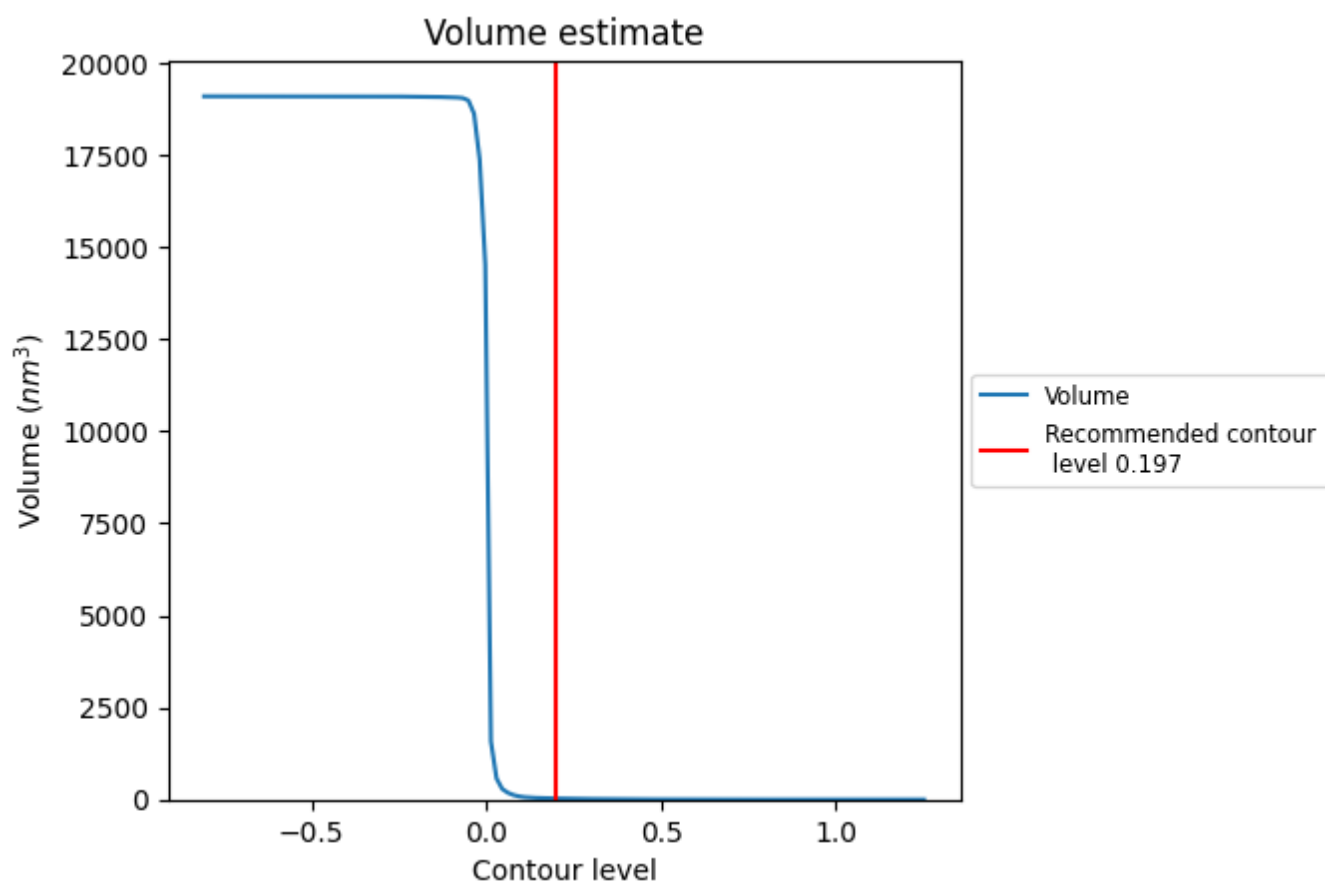
This section contains the results of statistical analysis of the map.

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.

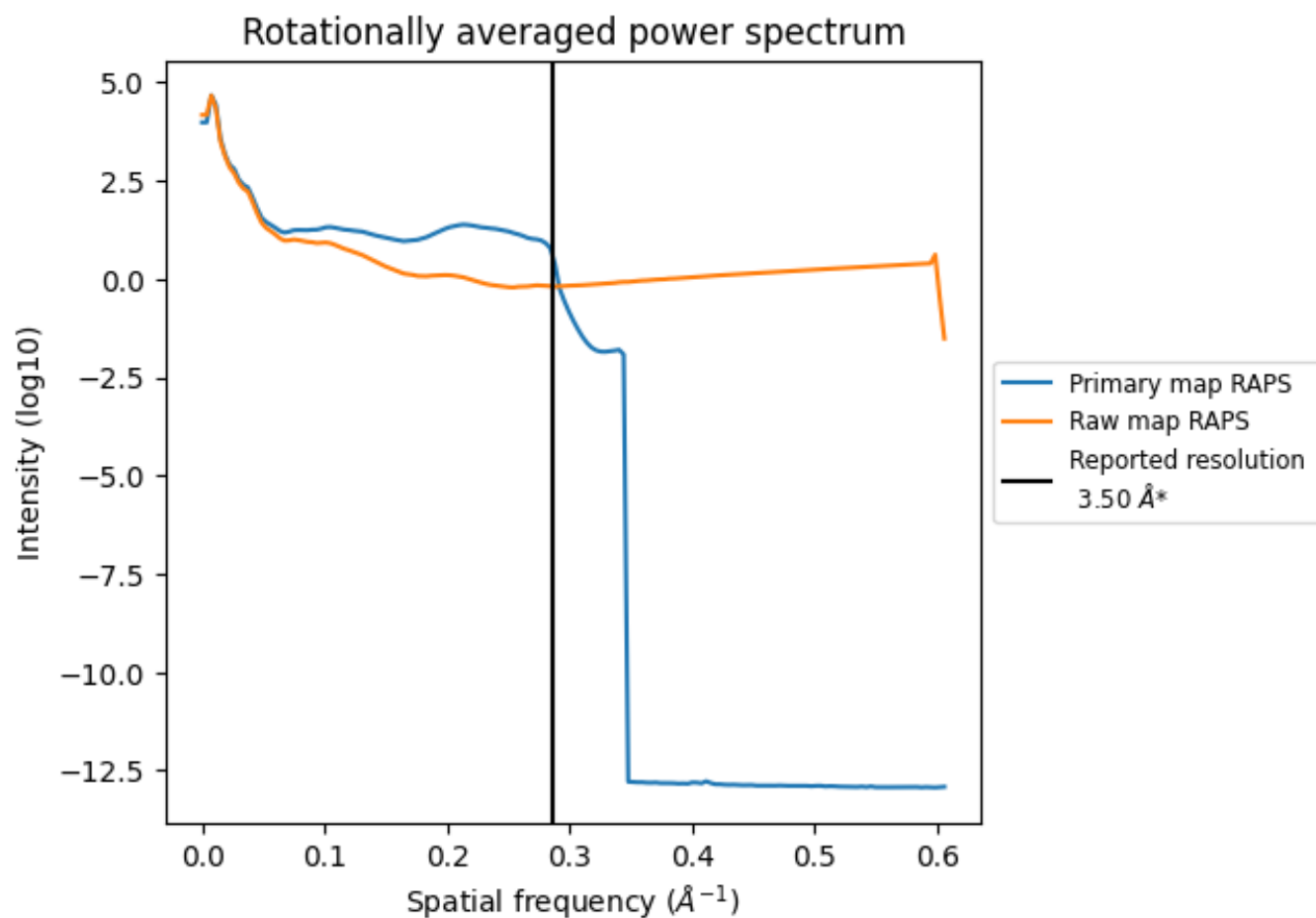
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 30 nm³; this corresponds to an approximate mass of 27 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 ⓘ

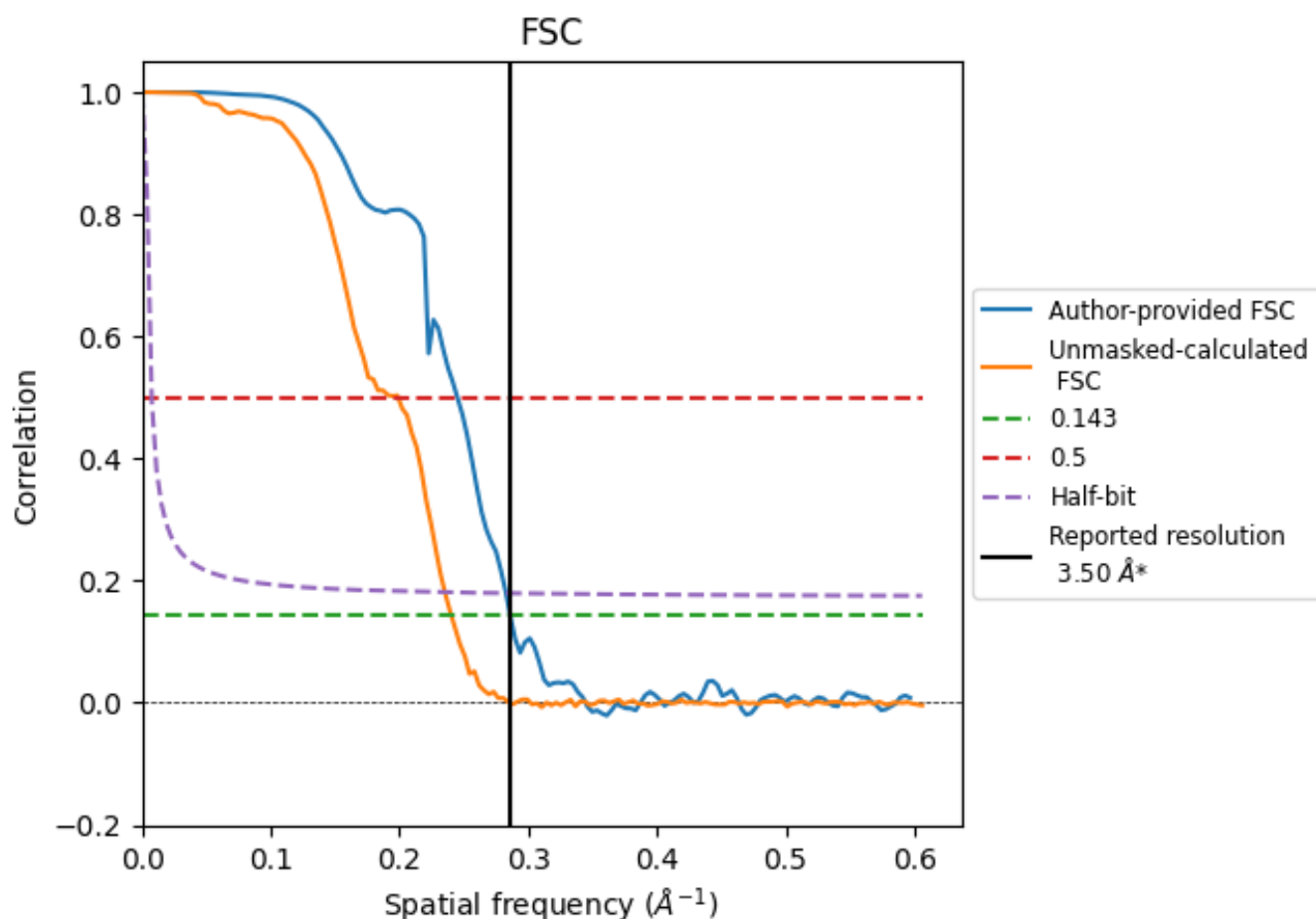


*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

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 0.286 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

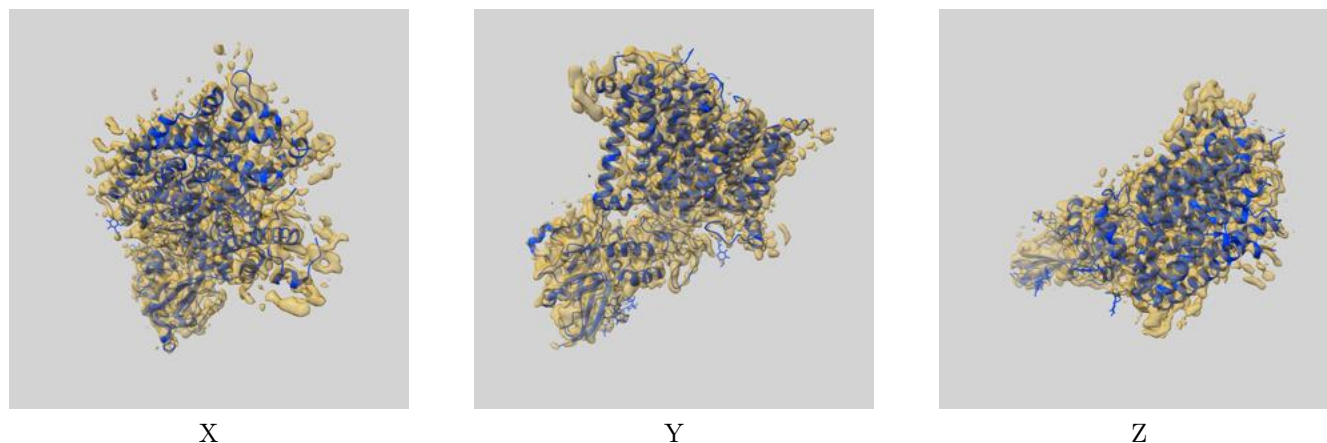
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	3.50	4.08	3.54
Unmasked-calculated*	4.16	5.03	4.25

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.16 differs from the reported value 3.5 by more than 10 %

9 Map-model fit [i](#)

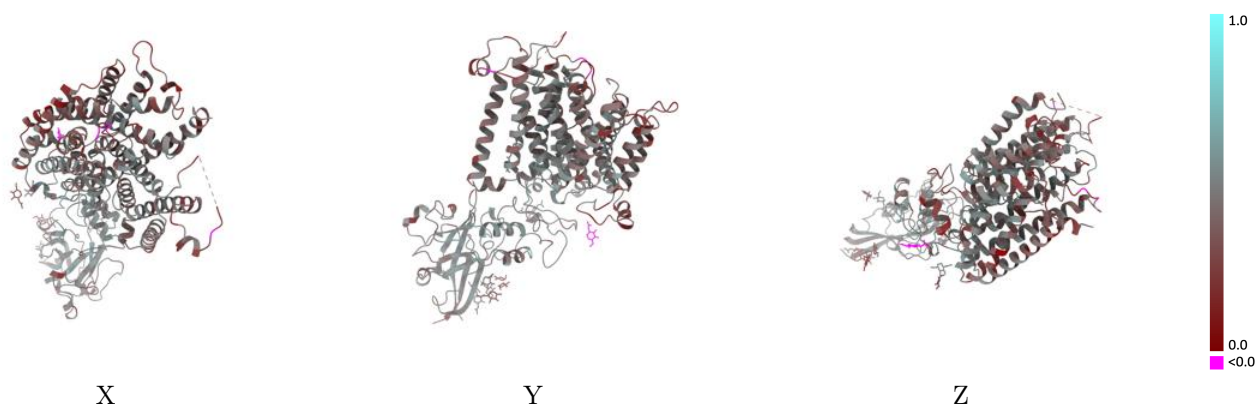
This section contains information regarding the fit between EMDB map EMD-50668 and PDB model 9FQT. Per-residue inclusion information can be found in section 3 on page 9.

9.1 Map-model overlay [i](#)



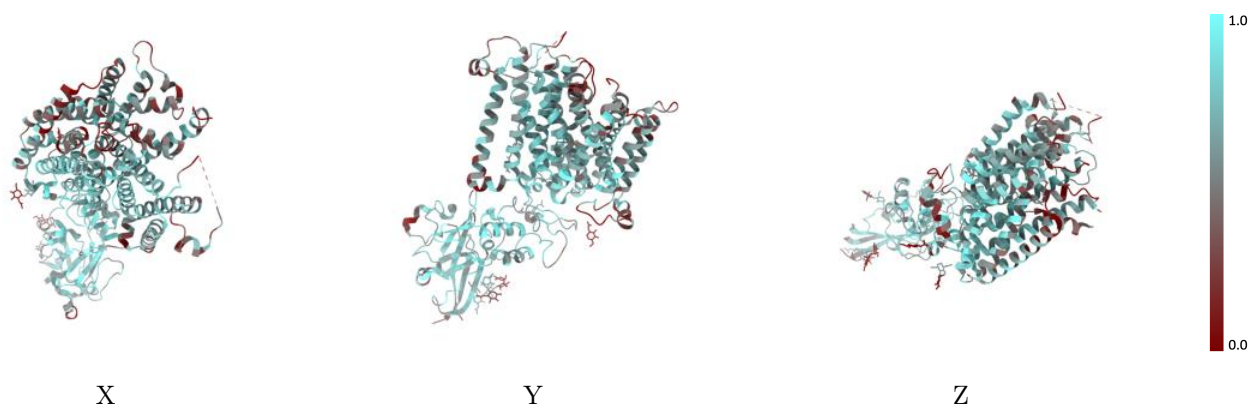
The images above show the 3D surface view of the map at the recommended contour level 0.197 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



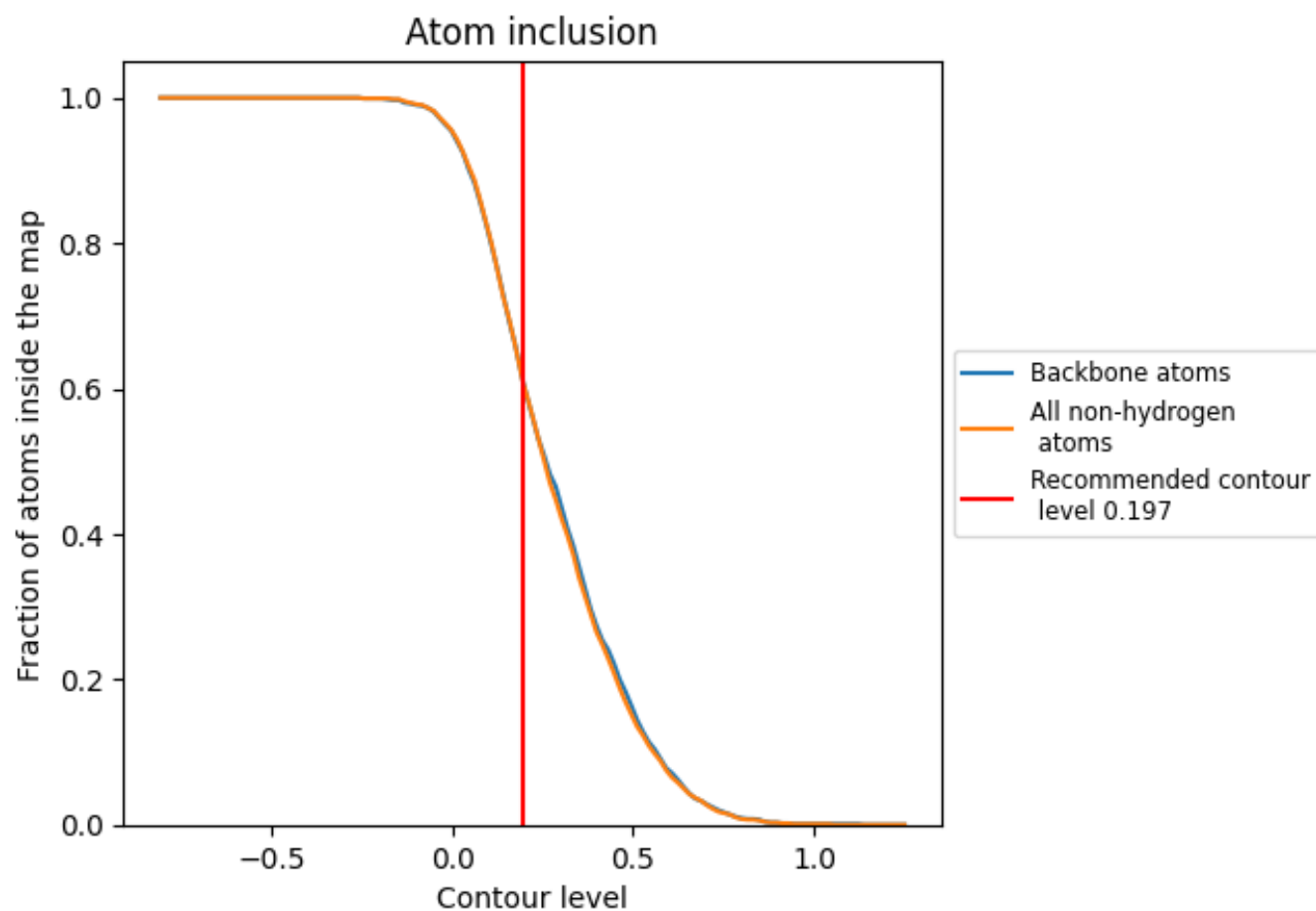
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.197).

9.4 Atom inclusion [i](#)



At the recommended contour level, 61% of all backbone atoms, 61% 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 (0.197) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6060	0.4260
A	0.5940	0.4080
B	0.7010	0.4710
C	0.2860	0.3930
D	0.3500	0.3140
I	0.4640	0.4100

