



# wwPDB EM Validation Summary Report ⓘ

Mar 23, 2026 – 11:09 PM UTC

PDB ID : 8D21 / pdb\_00008d21  
EMDB ID : EMD-27139  
Title : Cryo-EM structure of the VRC321 clinical trial, vaccine-elicited, human antibody 1B06 in complex with a stabilized NC99 HA trimer  
Authors : Gorman, J.; Kwong, P.D.  
Deposited on : 2022-05-27  
Resolution : 3.96 Å(reported)  
Based on initial model : 7MFG

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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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

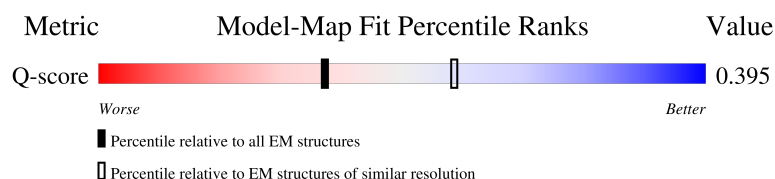
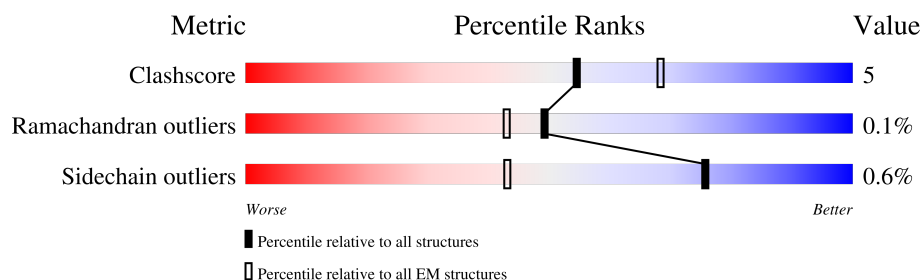
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.96 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	7646 ( 3.46 - 4.46 )

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion  $< 40\%$ ). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	B	222	 5% 71% 7% 22%
1	C	222	 5% 68% 10% 22%
1	G	222	 5% 72% 6% 22%
2	A	326	 85% 13% .

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Mol	Chain	Length	Quality of chain
2	D	326	
2	I	326	
3	E	123	
3	H	123	
3	J	123	
4	F	108	
4	K	108	
4	L	108	
5	M	3	
5	N	3	
5	O	3	

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 17541 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 Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	173	Total	C	N	O	S	0	0
			1396	875	239	274	8		
1	C	173	Total	C	N	O	S	0	0
			1396	875	239	274	8		
1	G	173	Total	C	N	O	S	0	0
			1396	875	239	274	8		

There are 141 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	47	CYS	GLY	conflict	UNP Q289M7
B	177	SER	-	expression tag	UNP Q289M7
B	178	GLY	-	expression tag	UNP Q289M7
B	179	ARG	-	expression tag	UNP Q289M7
B	180	LEU	-	expression tag	UNP Q289M7
B	181	VAL	-	expression tag	UNP Q289M7
B	182	PRO	-	expression tag	UNP Q289M7
B	183	ARG	-	expression tag	UNP Q289M7
B	184	GLY	-	expression tag	UNP Q289M7
B	185	SER	-	expression tag	UNP Q289M7
B	186	PRO	-	expression tag	UNP Q289M7
B	187	GLY	-	expression tag	UNP Q289M7
B	188	SER	-	expression tag	UNP Q289M7
B	189	GLY	-	expression tag	UNP Q289M7
B	190	TYR	-	expression tag	UNP Q289M7
B	191	ILE	-	expression tag	UNP Q289M7
B	192	PRO	-	expression tag	UNP Q289M7
B	193	GLU	-	expression tag	UNP Q289M7
B	194	ALA	-	expression tag	UNP Q289M7
B	195	PRO	-	expression tag	UNP Q289M7
B	196	ARG	-	expression tag	UNP Q289M7
B	197	ASP	-	expression tag	UNP Q289M7
B	198	GLY	-	expression tag	UNP Q289M7
B	199	GLN	-	expression tag	UNP Q289M7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	200	ALA	-	expression tag	UNP Q289M7
B	201	TYR	-	expression tag	UNP Q289M7
B	202	VAL	-	expression tag	UNP Q289M7
B	203	ARG	-	expression tag	UNP Q289M7
B	204	LYS	-	expression tag	UNP Q289M7
B	205	ASP	-	expression tag	UNP Q289M7
B	206	GLY	-	expression tag	UNP Q289M7
B	207	GLU	-	expression tag	UNP Q289M7
B	208	TRP	-	expression tag	UNP Q289M7
B	209	VAL	-	expression tag	UNP Q289M7
B	210	LEU	-	expression tag	UNP Q289M7
B	211	LEU	-	expression tag	UNP Q289M7
B	212	SER	-	expression tag	UNP Q289M7
B	213	THR	-	expression tag	UNP Q289M7
B	214	PHE	-	expression tag	UNP Q289M7
B	215	LEU	-	expression tag	UNP Q289M7
B	216	GLY	-	expression tag	UNP Q289M7
B	217	HIS	-	expression tag	UNP Q289M7
B	218	HIS	-	expression tag	UNP Q289M7
B	219	HIS	-	expression tag	UNP Q289M7
B	220	HIS	-	expression tag	UNP Q289M7
B	221	HIS	-	expression tag	UNP Q289M7
B	222	HIS	-	expression tag	UNP Q289M7
C	47	CYS	GLY	conflict	UNP Q289M7
C	177	SER	-	expression tag	UNP Q289M7
C	178	GLY	-	expression tag	UNP Q289M7
C	179	ARG	-	expression tag	UNP Q289M7
C	180	LEU	-	expression tag	UNP Q289M7
C	181	VAL	-	expression tag	UNP Q289M7
C	182	PRO	-	expression tag	UNP Q289M7
C	183	ARG	-	expression tag	UNP Q289M7
C	184	GLY	-	expression tag	UNP Q289M7
C	185	SER	-	expression tag	UNP Q289M7
C	186	PRO	-	expression tag	UNP Q289M7
C	187	GLY	-	expression tag	UNP Q289M7
C	188	SER	-	expression tag	UNP Q289M7
C	189	GLY	-	expression tag	UNP Q289M7
C	190	TYR	-	expression tag	UNP Q289M7
C	191	ILE	-	expression tag	UNP Q289M7
C	192	PRO	-	expression tag	UNP Q289M7
C	193	GLU	-	expression tag	UNP Q289M7
C	194	ALA	-	expression tag	UNP Q289M7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	195	PRO	-	expression tag	UNP Q289M7
C	196	ARG	-	expression tag	UNP Q289M7
C	197	ASP	-	expression tag	UNP Q289M7
C	198	GLY	-	expression tag	UNP Q289M7
C	199	GLN	-	expression tag	UNP Q289M7
C	200	ALA	-	expression tag	UNP Q289M7
C	201	TYR	-	expression tag	UNP Q289M7
C	202	VAL	-	expression tag	UNP Q289M7
C	203	ARG	-	expression tag	UNP Q289M7
C	204	LYS	-	expression tag	UNP Q289M7
C	205	ASP	-	expression tag	UNP Q289M7
C	206	GLY	-	expression tag	UNP Q289M7
C	207	GLU	-	expression tag	UNP Q289M7
C	208	TRP	-	expression tag	UNP Q289M7
C	209	VAL	-	expression tag	UNP Q289M7
C	210	LEU	-	expression tag	UNP Q289M7
C	211	LEU	-	expression tag	UNP Q289M7
C	212	SER	-	expression tag	UNP Q289M7
C	213	THR	-	expression tag	UNP Q289M7
C	214	PHE	-	expression tag	UNP Q289M7
C	215	LEU	-	expression tag	UNP Q289M7
C	216	GLY	-	expression tag	UNP Q289M7
C	217	HIS	-	expression tag	UNP Q289M7
C	218	HIS	-	expression tag	UNP Q289M7
C	219	HIS	-	expression tag	UNP Q289M7
C	220	HIS	-	expression tag	UNP Q289M7
C	221	HIS	-	expression tag	UNP Q289M7
C	222	HIS	-	expression tag	UNP Q289M7
G	47	CYS	GLY	conflict	UNP Q289M7
G	177	SER	-	expression tag	UNP Q289M7
G	178	GLY	-	expression tag	UNP Q289M7
G	179	ARG	-	expression tag	UNP Q289M7
G	180	LEU	-	expression tag	UNP Q289M7
G	181	VAL	-	expression tag	UNP Q289M7
G	182	PRO	-	expression tag	UNP Q289M7
G	183	ARG	-	expression tag	UNP Q289M7
G	184	GLY	-	expression tag	UNP Q289M7
G	185	SER	-	expression tag	UNP Q289M7
G	186	PRO	-	expression tag	UNP Q289M7
G	187	GLY	-	expression tag	UNP Q289M7
G	188	SER	-	expression tag	UNP Q289M7
G	189	GLY	-	expression tag	UNP Q289M7

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Chain	Residue	Modelled	Actual	Comment	Reference
G	190	TYR	-	expression tag	UNP Q289M7
G	191	ILE	-	expression tag	UNP Q289M7
G	192	PRO	-	expression tag	UNP Q289M7
G	193	GLU	-	expression tag	UNP Q289M7
G	194	ALA	-	expression tag	UNP Q289M7
G	195	PRO	-	expression tag	UNP Q289M7
G	196	ARG	-	expression tag	UNP Q289M7
G	197	ASP	-	expression tag	UNP Q289M7
G	198	GLY	-	expression tag	UNP Q289M7
G	199	GLN	-	expression tag	UNP Q289M7
G	200	ALA	-	expression tag	UNP Q289M7
G	201	TYR	-	expression tag	UNP Q289M7
G	202	VAL	-	expression tag	UNP Q289M7
G	203	ARG	-	expression tag	UNP Q289M7
G	204	LYS	-	expression tag	UNP Q289M7
G	205	ASP	-	expression tag	UNP Q289M7
G	206	GLY	-	expression tag	UNP Q289M7
G	207	GLU	-	expression tag	UNP Q289M7
G	208	TRP	-	expression tag	UNP Q289M7
G	209	VAL	-	expression tag	UNP Q289M7
G	210	LEU	-	expression tag	UNP Q289M7
G	211	LEU	-	expression tag	UNP Q289M7
G	212	SER	-	expression tag	UNP Q289M7
G	213	THR	-	expression tag	UNP Q289M7
G	214	PHE	-	expression tag	UNP Q289M7
G	215	LEU	-	expression tag	UNP Q289M7
G	216	GLY	-	expression tag	UNP Q289M7
G	217	HIS	-	expression tag	UNP Q289M7
G	218	HIS	-	expression tag	UNP Q289M7
G	219	HIS	-	expression tag	UNP Q289M7
G	220	HIS	-	expression tag	UNP Q289M7
G	221	HIS	-	expression tag	UNP Q289M7
G	222	HIS	-	expression tag	UNP Q289M7

- Molecule 2 is a protein called Hemagglutinin HA1 chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	323	Total	C	N	O	S	0	0
			2536	1604	437	483	12		
2	D	323	Total	C	N	O	S	0	0
			2536	1604	437	483	12		
2	I	323	Total	C	N	O	S	0	0
			2536	1604	437	483	12		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	CYS	LEU	conflict	UNP Q6WG00
D	30	CYS	LEU	conflict	UNP Q6WG00
I	30	CYS	LEU	conflict	UNP Q6WG00

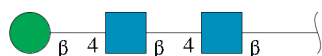
- Molecule 3 is a protein called 1B06 Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	H	123	Total	C	N	O	S	0	0
			950	596	171	180	3		
3	E	123	Total	C	N	O	S	0	0
			950	596	171	180	3		
3	J	123	Total	C	N	O	S	0	0
			950	596	171	180	3		

- Molecule 4 is a protein called 1B06 Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	L	108	Total	C	N	O	S	0	0
			828	520	141	164	3		
4	F	108	Total	C	N	O	S	0	0
			828	520	141	164	3		
4	K	108	Total	C	N	O	S	0	0
			828	520	141	164	3		

- Molecule 5 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
5	M	3	Total	C	N	O		0	0
			39	22	2	15			
5	N	3	Total	C	N	O		0	0
			39	22	2	15			
5	O	3	Total	C	N	O		0	0
			39	22	2	15			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).





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	A	1	Total	C	N	O	0
			14	8	1	5	
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	A	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	
6	D	1	Total	C	N	O	0
			14	8	1	5	

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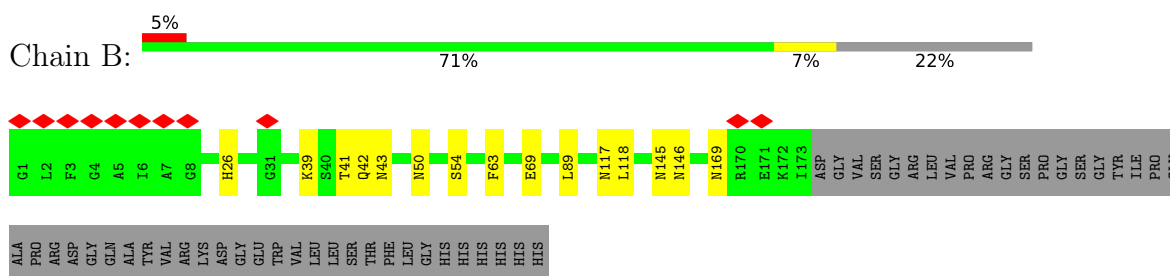
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Mol	Chain	Residues	Atoms				AltConf
6	I	1	Total	C	N	O	0
			14	8	1	5	
6	I	1	Total	C	N	O	0
			14	8	1	5	
6	I	1	Total	C	N	O	0
			14	8	1	5	
6	I	1	Total	C	N	O	0
			14	8	1	5	
6	I	1	Total	C	N	O	0
			14	8	1	5	
6	I	1	Total	C	N	O	0
			14	8	1	5	

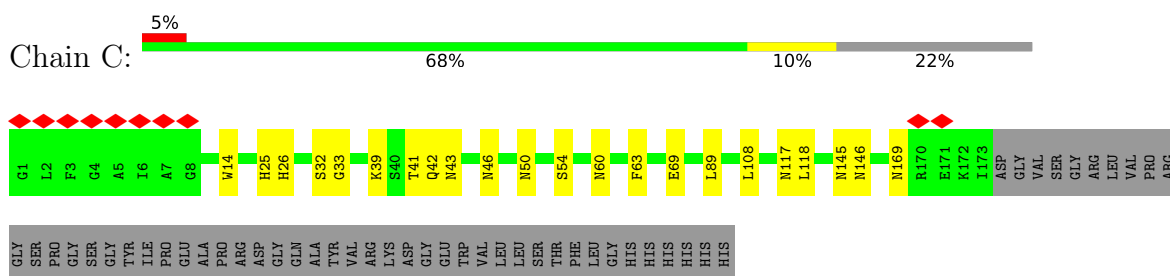
### 3 Residue-property plots [i](#)

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

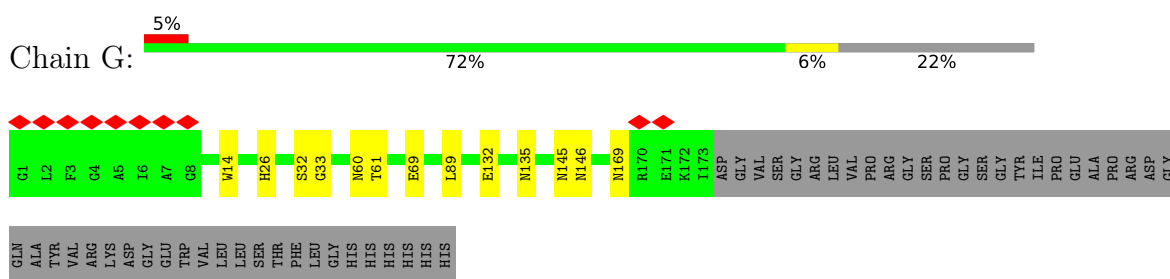
- Molecule 1: Hemagglutinin HA2 chain



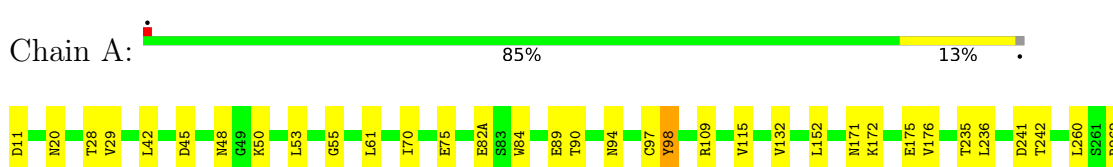
- Molecule 1: Hemagglutinin HA2 chain

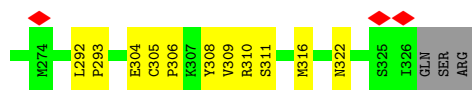


- Molecule 1: Hemagglutinin HA2 chain



- Molecule 2: Hemagglutinin HA1 chain





- Molecule 2: Hemagglutinin HA1 chain

Chain D: 87% 13%



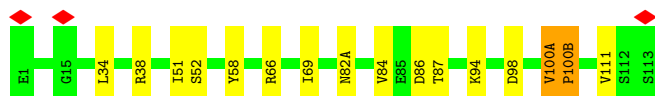
- Molecule 2: Hemagglutinin HA1 chain

Chain I: 85% 14%



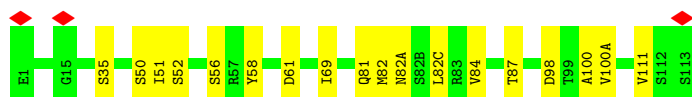
- Molecule 3: 1B06 Heavy Chain

Chain H: 87% 11%



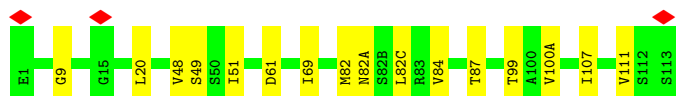
- Molecule 3: 1B06 Heavy Chain

Chain E: 85% 15%



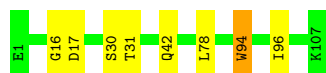
- Molecule 3: 1B06 Heavy Chain

Chain J: 87% 13%



- Molecule 4: 1B06 Light Chain

Chain L:  93% 6%



- Molecule 4: 1B06 Light Chain

Chain F:  92% 8%



- Molecule 4: 1B06 Light Chain

Chain K:  94% 6%



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

Chain M:  33% 67%



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

Chain N:  33% 67%



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

Chain O:  33% 67%



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C3	Depositor
Number of particles used	73396	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 ( $e^-/\text{\AA}^2$ )	70.72	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.228	Depositor
Minimum map value	-0.909	Depositor
Average map value	0.014	Depositor
Map value standard deviation	0.086	Depositor
Recommended contour level	0.6	Depositor
Map size (Å)	324.75, 324.75, 324.75	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.0825, 1.0825, 1.0825	Depositor

## 5 Model quality

### 5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	B	0.16	0/1423	0.34	0/1912
1	C	0.15	0/1423	0.32	0/1912
1	G	0.15	0/1423	0.32	0/1912
2	A	0.16	0/2604	0.33	0/3549
2	D	0.17	0/2604	0.37	1/3549 (0.0%)
2	I	0.16	0/2604	0.35	0/3549
3	E	0.15	0/970	0.34	0/1310
3	H	0.15	0/970	0.35	0/1310
3	J	0.15	0/970	0.32	0/1310
4	F	0.16	0/847	0.35	0/1153
4	K	0.18	0/847	0.37	0/1153
4	L	0.19	0/847	0.42	0/1153
All	All	0.16	0/17532	0.35	1/23772 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	H	0	1
4	K	0	1
4	L	0	1
All	All	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	306	PRO	CA-N-CD	-7.40	101.63	112.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	H	100(A)	VAL	Peptide
4	K	94	TRP	Peptide
4	L	94	TRP	Peptide

## 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	B	1396	0	1328	10	0
1	C	1396	0	1328	17	0
1	G	1396	0	1328	10	0
2	A	2536	0	2440	32	0
2	D	2536	0	2440	31	0
2	I	2536	0	2440	30	0
3	E	950	0	933	13	0
3	H	950	0	933	10	0
3	J	950	0	933	12	0
4	F	828	0	805	5	0
4	K	828	0	805	4	0
4	L	828	0	805	6	0
5	M	39	0	34	1	0
5	N	39	0	34	2	0
5	O	39	0	34	2	0
6	A	98	0	91	3	0
6	D	98	0	91	3	0
6	I	98	0	91	2	0
All	All	17541	0	16893	174	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 5.

The worst 5 of 174 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)
4:L:94:TRP:HE1	4:L:96:ILE:HD11	1.55	0.71
1:B:39:LYS:O	1:B:43:ASN:ND2	2.23	0.71
3:J:51:ILE:HD13	3:J:69:ILE:HG21	1.72	0.70
2:A:11:ASP:N	3:H:58:TYR:HH	1.91	0.69
2:D:11:ASP:N	3:E:58:TYR:HH	1.91	0.69

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	171/222 (77%)	161 (94%)	10 (6%)	0	100	100
1	C	171/222 (77%)	162 (95%)	9 (5%)	0	100	100
1	G	171/222 (77%)	163 (95%)	8 (5%)	0	100	100
2	A	321/326 (98%)	299 (93%)	22 (7%)	0	100	100
2	D	321/326 (98%)	306 (95%)	15 (5%)	0	100	100
2	I	321/326 (98%)	304 (95%)	17 (5%)	0	100	100
3	E	121/123 (98%)	116 (96%)	4 (3%)	1 (1%)	16	51
3	H	121/123 (98%)	113 (93%)	7 (6%)	1 (1%)	16	51
3	J	121/123 (98%)	116 (96%)	5 (4%)	0	100	100
4	F	106/108 (98%)	93 (88%)	13 (12%)	0	100	100
4	K	106/108 (98%)	96 (91%)	10 (9%)	0	100	100
4	L	106/108 (98%)	93 (88%)	13 (12%)	0	100	100
All	All	2157/2337 (92%)	2022 (94%)	133 (6%)	2 (0%)	49	81

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	H	100(B)	PRO
3	E	100	ALA

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	150/189 (79%)	149 (99%)	1 (1%)	76	79
1	C	150/189 (79%)	148 (99%)	2 (1%)	61	72
1	G	150/189 (79%)	149 (99%)	1 (1%)	76	79
2	A	284/287 (99%)	280 (99%)	4 (1%)	59	71
2	D	284/287 (99%)	283 (100%)	1 (0%)	84	84
2	I	284/287 (99%)	283 (100%)	1 (0%)	84	84
3	E	104/104 (100%)	104 (100%)	0	100	100
3	H	104/104 (100%)	103 (99%)	1 (1%)	68	76
3	J	104/104 (100%)	104 (100%)	0	100	100
4	F	91/91 (100%)	90 (99%)	1 (1%)	65	74
4	K	91/91 (100%)	91 (100%)	0	100	100
4	L	91/91 (100%)	91 (100%)	0	100	100
All	All	1887/2013 (94%)	1875 (99%)	12 (1%)	76	81

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	169	ASN
2	D	203	SER
2	I	136	SER
4	F	88	CYS
2	A	98	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	190	ASN
2	D	226	GLN
2	I	226	GLN
1	G	169	ASN
2	I	183	HIS

### 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 ⓘ

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	$\# Z  > 2$
5	NAG	M	1	1,5	14,14,15	0.26	0	17,19,21	0.47	0
5	NAG	M	2	5	14,14,15	0.19	0	17,19,21	0.66	0
5	BMA	M	3	5	11,11,12	0.53	0	15,15,17	0.68	0
5	NAG	N	1	1,5	14,14,15	0.24	0	17,19,21	0.46	0
5	NAG	N	2	5	14,14,15	0.20	0	17,19,21	0.65	0
5	BMA	N	3	5	11,11,12	0.55	0	15,15,17	0.68	0
5	NAG	O	1	1,5	14,14,15	0.27	0	17,19,21	0.47	0
5	NAG	O	2	5	14,14,15	0.19	0	17,19,21	0.66	0
5	BMA	O	3	5	11,11,12	0.55	0	15,15,17	0.68	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	M	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	M	2	5	-	4/6/23/26	0/1/1/1
5	BMA	M	3	5	-	1/2/19/22	0/1/1/1
5	NAG	N	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	N	2	5	-	4/6/23/26	0/1/1/1
5	BMA	N	3	5	-	1/2/19/22	0/1/1/1
5	NAG	O	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	O	2	5	-	4/6/23/26	0/1/1/1
5	BMA	O	3	5	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 27 torsion outliers are listed below:

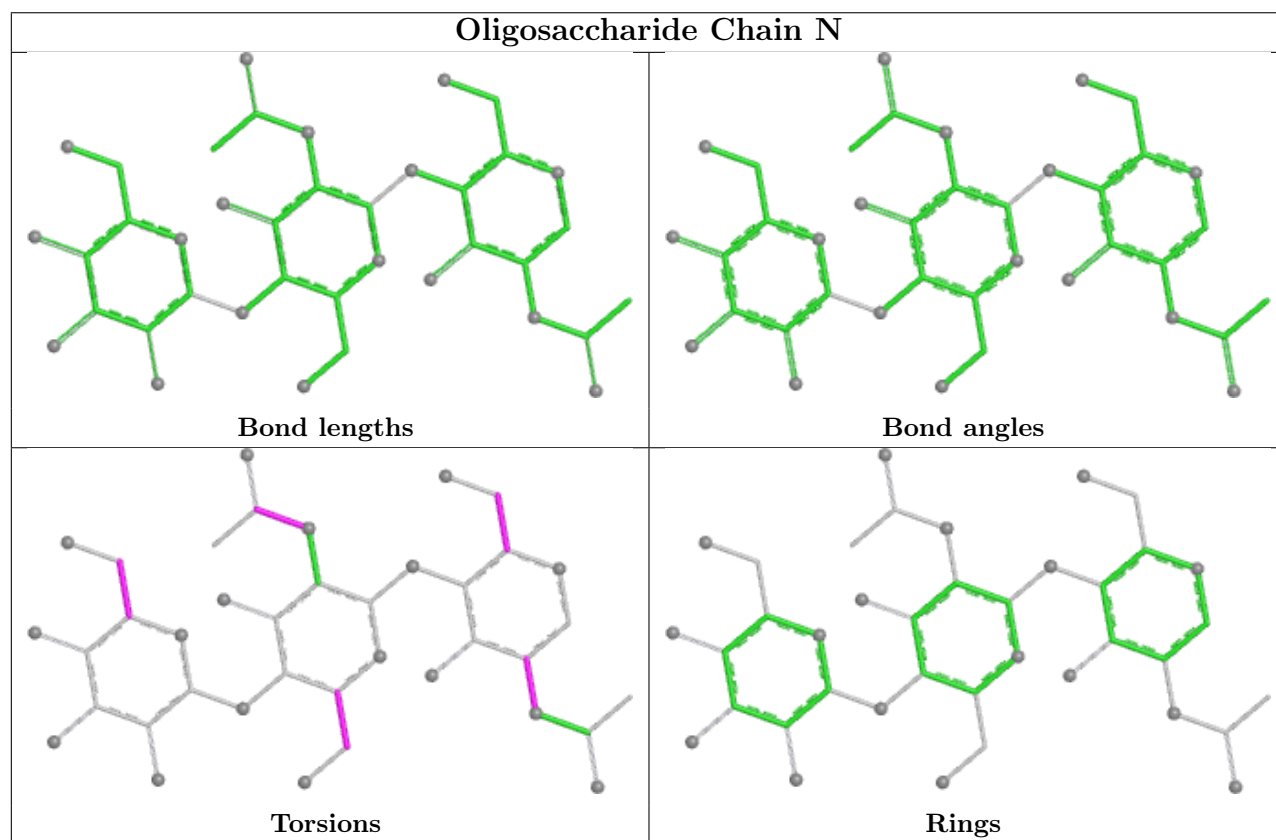
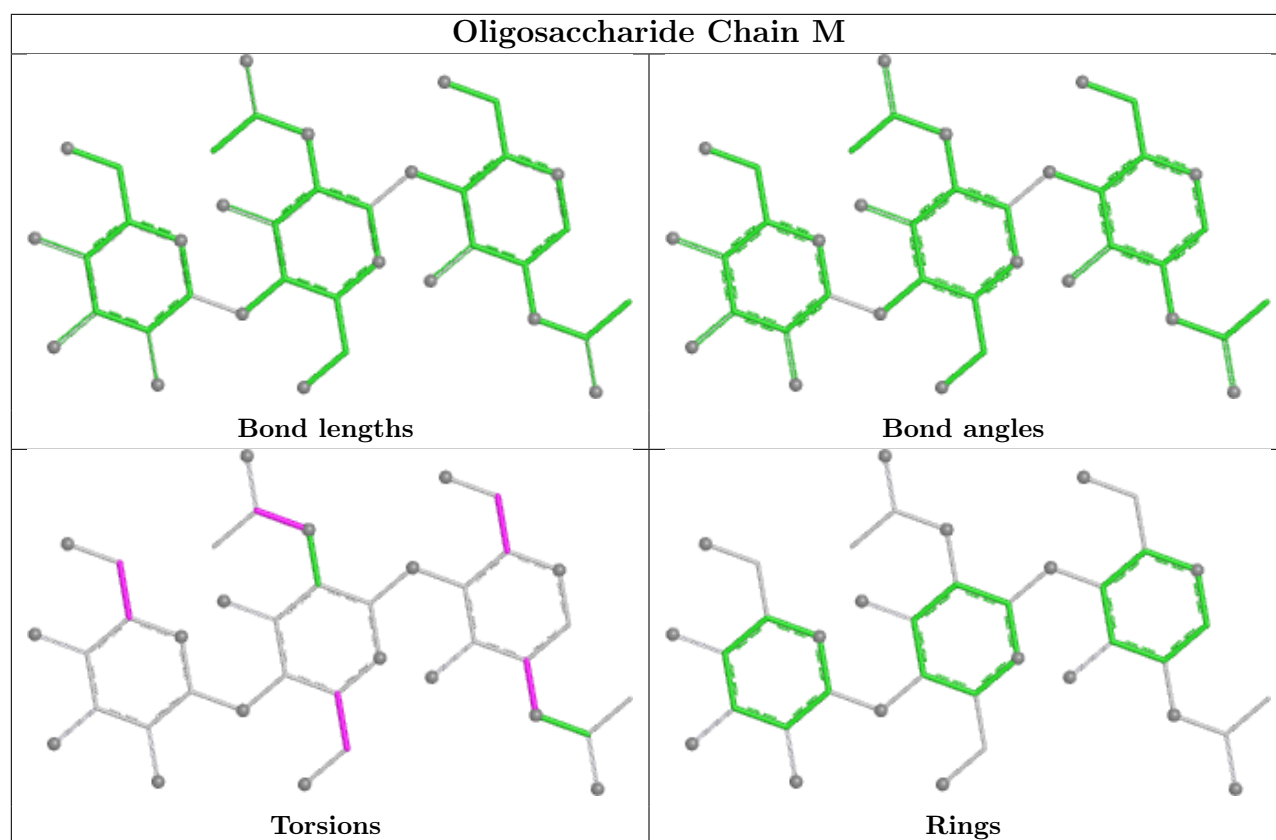
Mol	Chain	Res	Type	Atoms
5	O	2	NAG	O5-C5-C6-O6
5	M	2	NAG	O5-C5-C6-O6
5	N	1	NAG	O5-C5-C6-O6
5	N	2	NAG	O5-C5-C6-O6
5	O	1	NAG	O5-C5-C6-O6

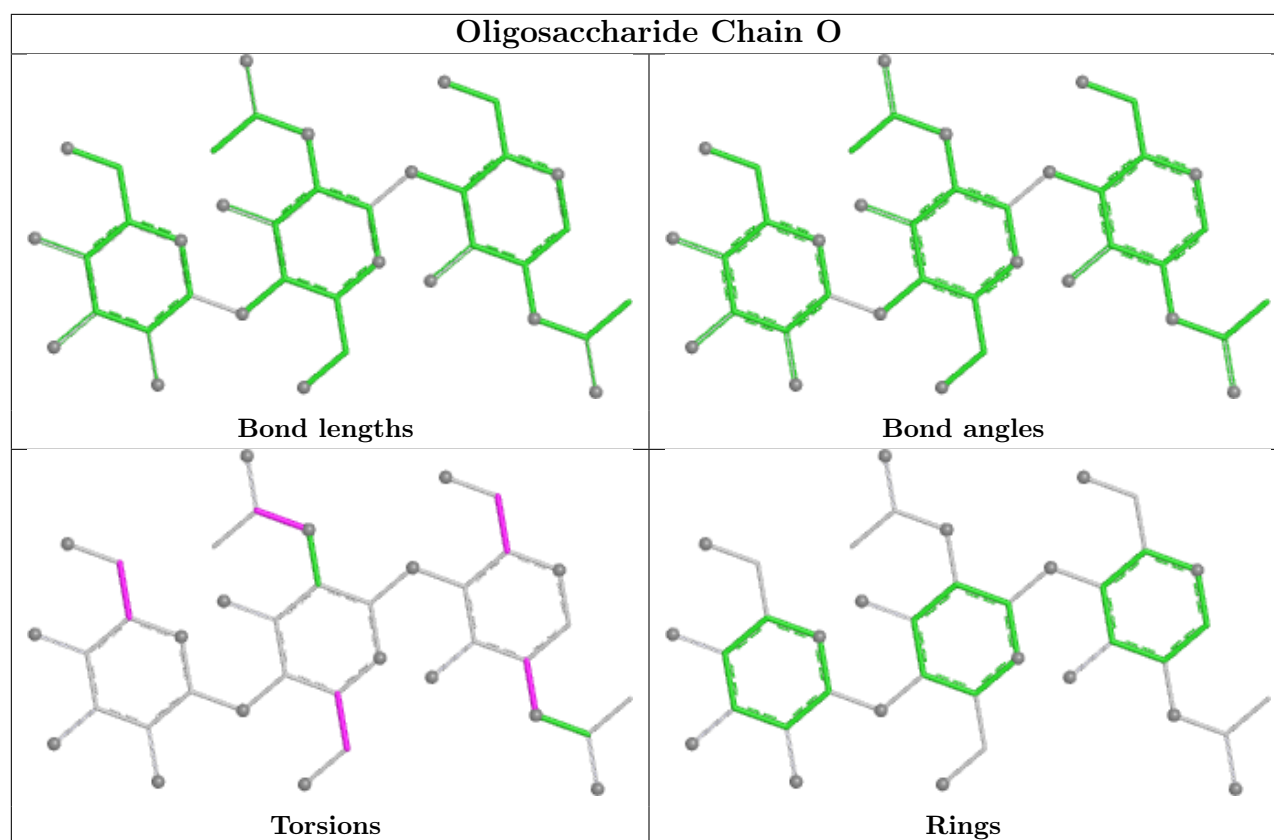
There are no ring outliers.

6 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	N	2	NAG	2	0
5	M	2	NAG	1	0
5	O	2	NAG	2	0
5	O	1	NAG	2	0
5	M	1	NAG	1	0
5	N	1	NAG	2	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](#)

21 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	I	404	2	14,14,15	0.20	0	17,19,21	0.43	0
6	NAG	A	405	2	14,14,15	0.20	0	17,19,21	0.44	0
6	NAG	D	403	2	14,14,15	0.35	0	17,19,21	0.49	0
6	NAG	I	406	2	14,14,15	0.19	0	17,19,21	0.47	0
6	NAG	D	404	2	14,14,15	0.21	0	17,19,21	0.44	0
6	NAG	I	401	2	14,14,15	0.35	0	17,19,21	0.51	0
6	NAG	I	402	2	14,14,15	0.24	0	17,19,21	1.19	1 (5%)
6	NAG	A	407	2	14,14,15	0.19	0	17,19,21	0.41	0
6	NAG	D	406	2	14,14,15	0.21	0	17,19,21	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	NAG	I	407	2	14,14,15	0.20	0	17,19,21	0.42	0
6	NAG	D	405	2	14,14,15	0.19	0	17,19,21	0.44	0
6	NAG	D	402	2	14,14,15	0.25	0	17,19,21	1.19	1 (5%)
6	NAG	D	401	2	14,14,15	0.34	0	17,19,21	0.50	0
6	NAG	A	402	2	14,14,15	0.25	0	17,19,21	1.20	1 (5%)
6	NAG	A	401	2	14,14,15	0.34	0	17,19,21	0.48	0
6	NAG	A	403	2	14,14,15	0.30	0	17,19,21	0.51	0
6	NAG	D	407	2	14,14,15	0.18	0	17,19,21	0.42	0
6	NAG	I	405	2	14,14,15	0.20	0	17,19,21	0.44	0
6	NAG	I	403	2	14,14,15	0.32	0	17,19,21	0.51	0
6	NAG	A	404	2	14,14,15	0.20	0	17,19,21	0.43	0
6	NAG	A	406	2	14,14,15	0.20	0	17,19,21	0.48	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	I	404	2	-	0/6/23/26	0/1/1/1
6	NAG	A	405	2	-	0/6/23/26	0/1/1/1
6	NAG	D	403	2	-	1/6/23/26	0/1/1/1
6	NAG	I	406	2	-	2/6/23/26	0/1/1/1
6	NAG	D	404	2	-	1/6/23/26	0/1/1/1
6	NAG	I	401	2	-	2/6/23/26	0/1/1/1
6	NAG	I	402	2	-	6/6/23/26	0/1/1/1
6	NAG	A	407	2	-	2/6/23/26	0/1/1/1
6	NAG	D	406	2	-	2/6/23/26	0/1/1/1
6	NAG	I	407	2	-	2/6/23/26	0/1/1/1
6	NAG	D	405	2	-	0/6/23/26	0/1/1/1
6	NAG	D	402	2	-	6/6/23/26	0/1/1/1
6	NAG	D	401	2	-	2/6/23/26	0/1/1/1
6	NAG	A	402	2	-	6/6/23/26	0/1/1/1
6	NAG	A	401	2	-	2/6/23/26	0/1/1/1
6	NAG	A	403	2	-	1/6/23/26	0/1/1/1
6	NAG	D	407	2	-	2/6/23/26	0/1/1/1
6	NAG	I	405	2	-	0/6/23/26	0/1/1/1
6	NAG	I	403	2	-	1/6/23/26	0/1/1/1
6	NAG	A	404	2	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	A	406	2	-	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(°)
6	A	402	NAG	C2-N2-C7	3.84	128.04	122.90
6	I	402	NAG	C2-N2-C7	3.82	128.03	122.90
6	D	402	NAG	C2-N2-C7	3.82	128.02	122.90

There are no chirality outliers.

5 of 41 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	402	NAG	O5-C5-C6-O6
6	I	402	NAG	O5-C5-C6-O6
6	D	402	NAG	O5-C5-C6-O6
6	A	402	NAG	C4-C5-C6-O6
6	D	402	NAG	C4-C5-C6-O6

There are no ring outliers.

6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	403	NAG	1	0
6	I	402	NAG	1	0
6	D	402	NAG	2	0
6	A	402	NAG	2	0
6	A	403	NAG	1	0
6	I	403	NAG	1	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-27139. These allow visual inspection of the internal detail of the map and identification of artifacts.

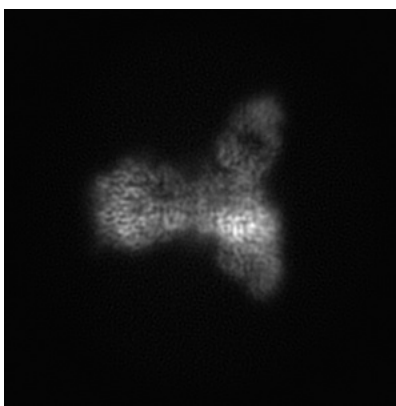
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

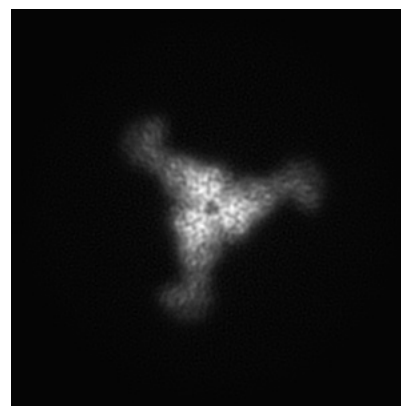
#### 6.1.1 Primary map



X

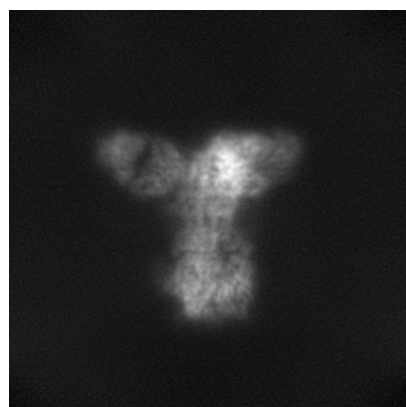


Y

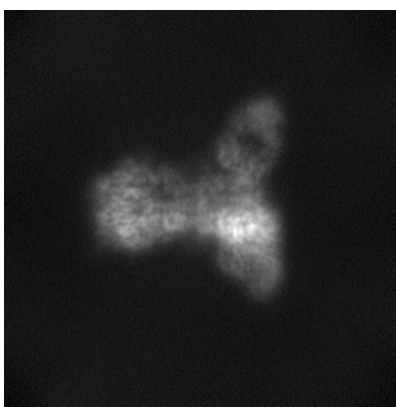


Z

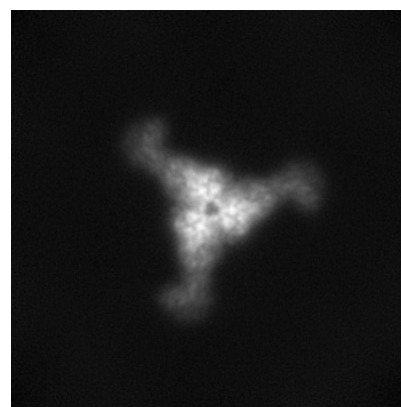
#### 6.1.2 Raw map



X



Y

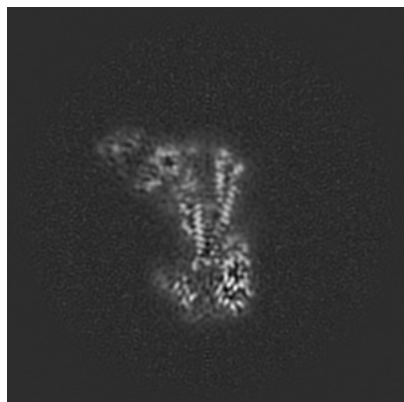


Z

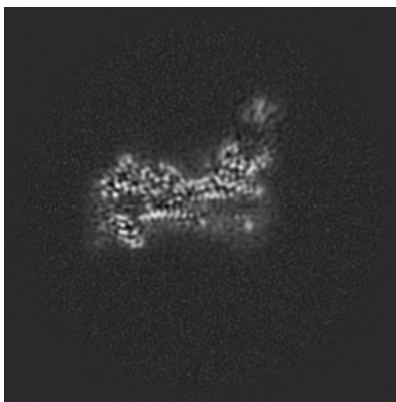
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

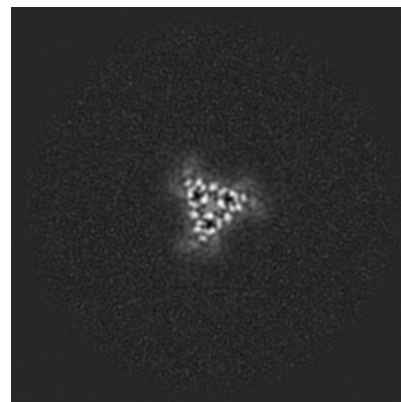
### 6.2.1 Primary map



X Index: 150

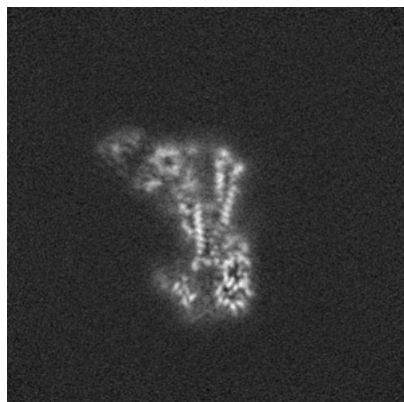


Y Index: 150

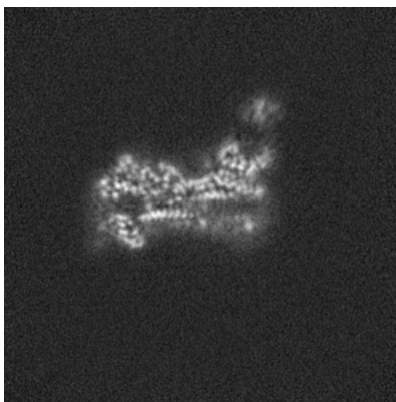


Z Index: 150

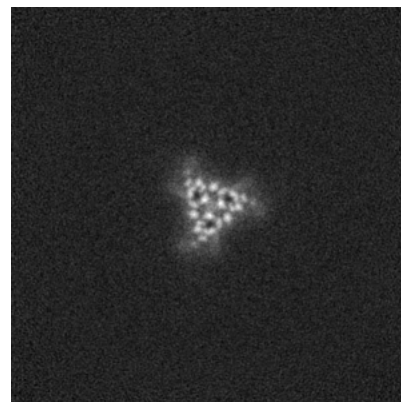
### 6.2.2 Raw map



X Index: 150



Y Index: 150

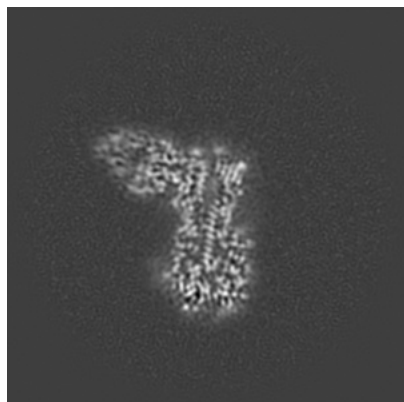


Z Index: 150

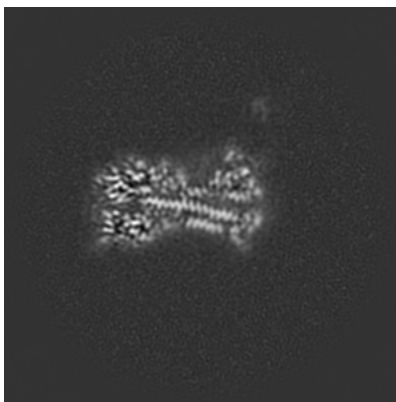
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

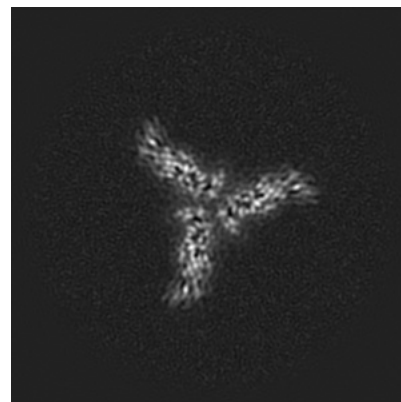
### 6.3.1 Primary map



X Index: 142

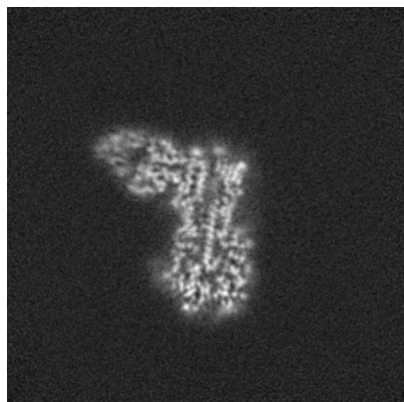


Y Index: 144

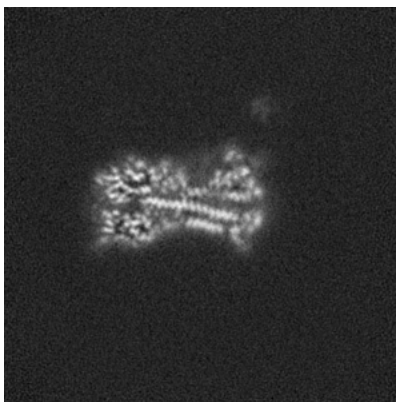


Z Index: 173

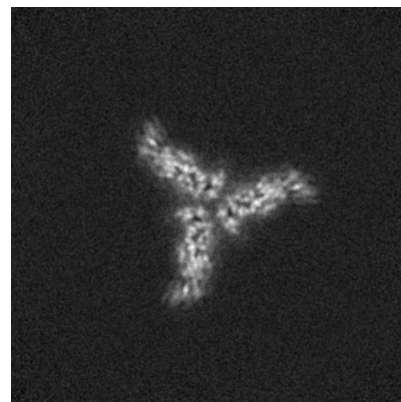
### 6.3.2 Raw map



X Index: 142



Y Index: 144

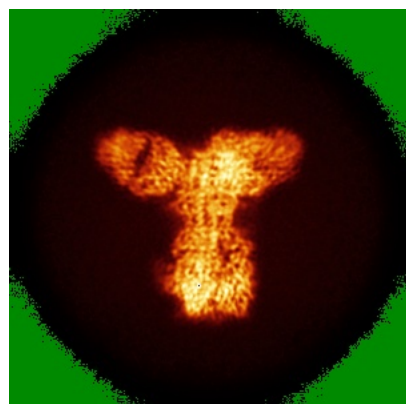


Z Index: 173

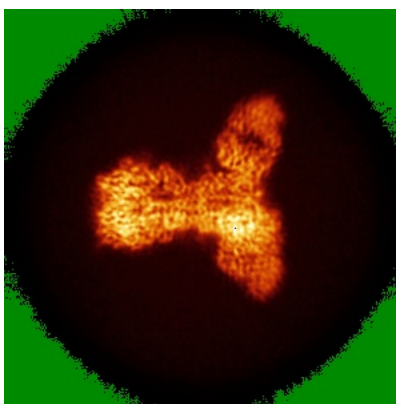
The images above show the largest variance slices of the map in three orthogonal directions.

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

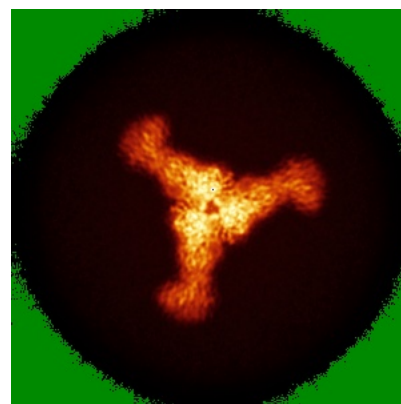
### 6.4.1 Primary map



X

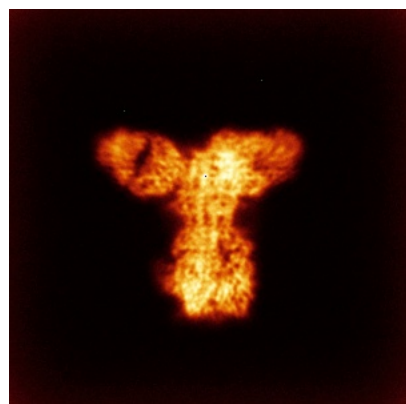


Y

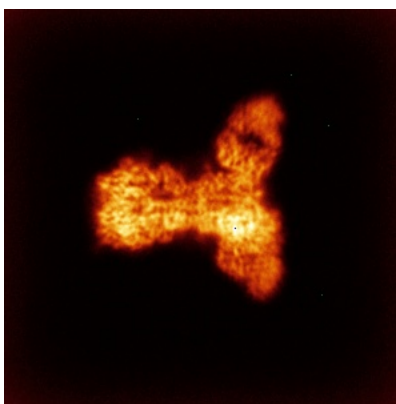


Z

### 6.4.2 Raw map



X



Y

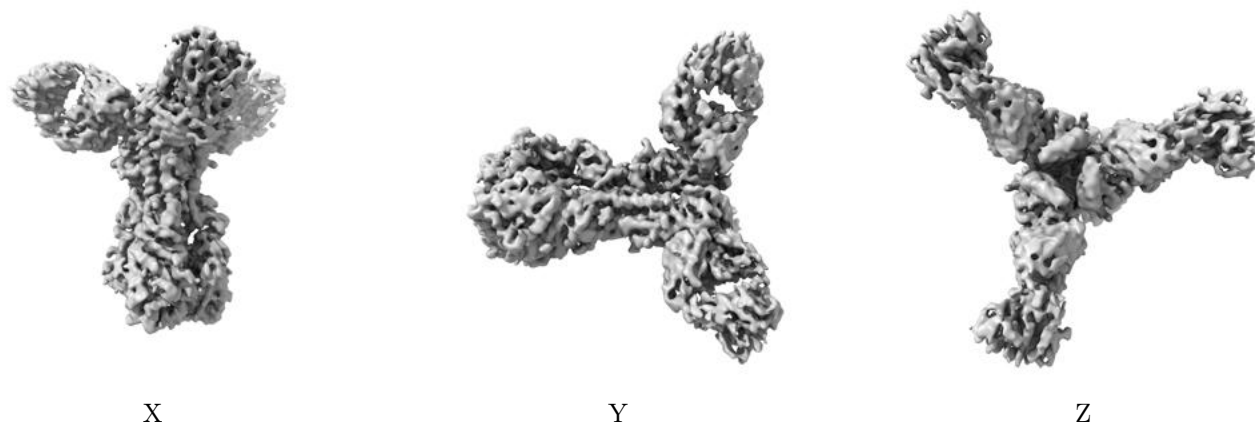


Z

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.

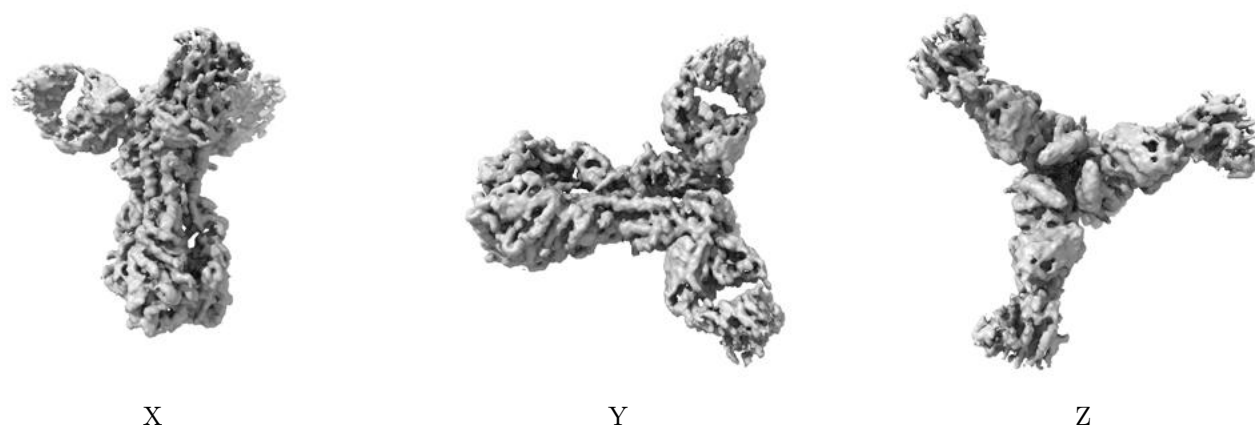
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.6. 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.



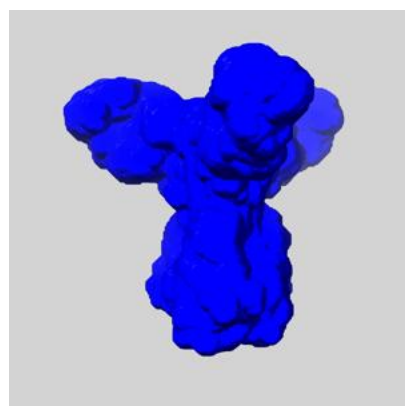
## 6.6 Mask visualisation [i](#)

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

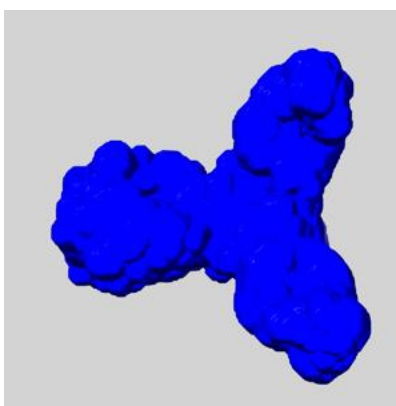
A mask typically either:

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

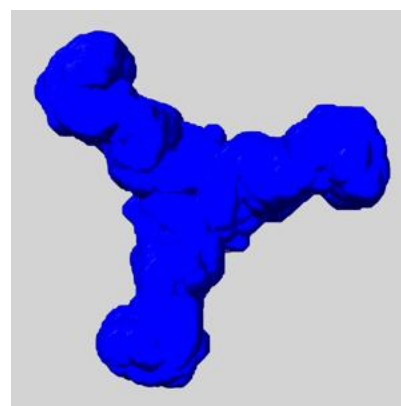
### 6.6.1 emd\_27139\_msk\_1.map [i](#)



X



Y

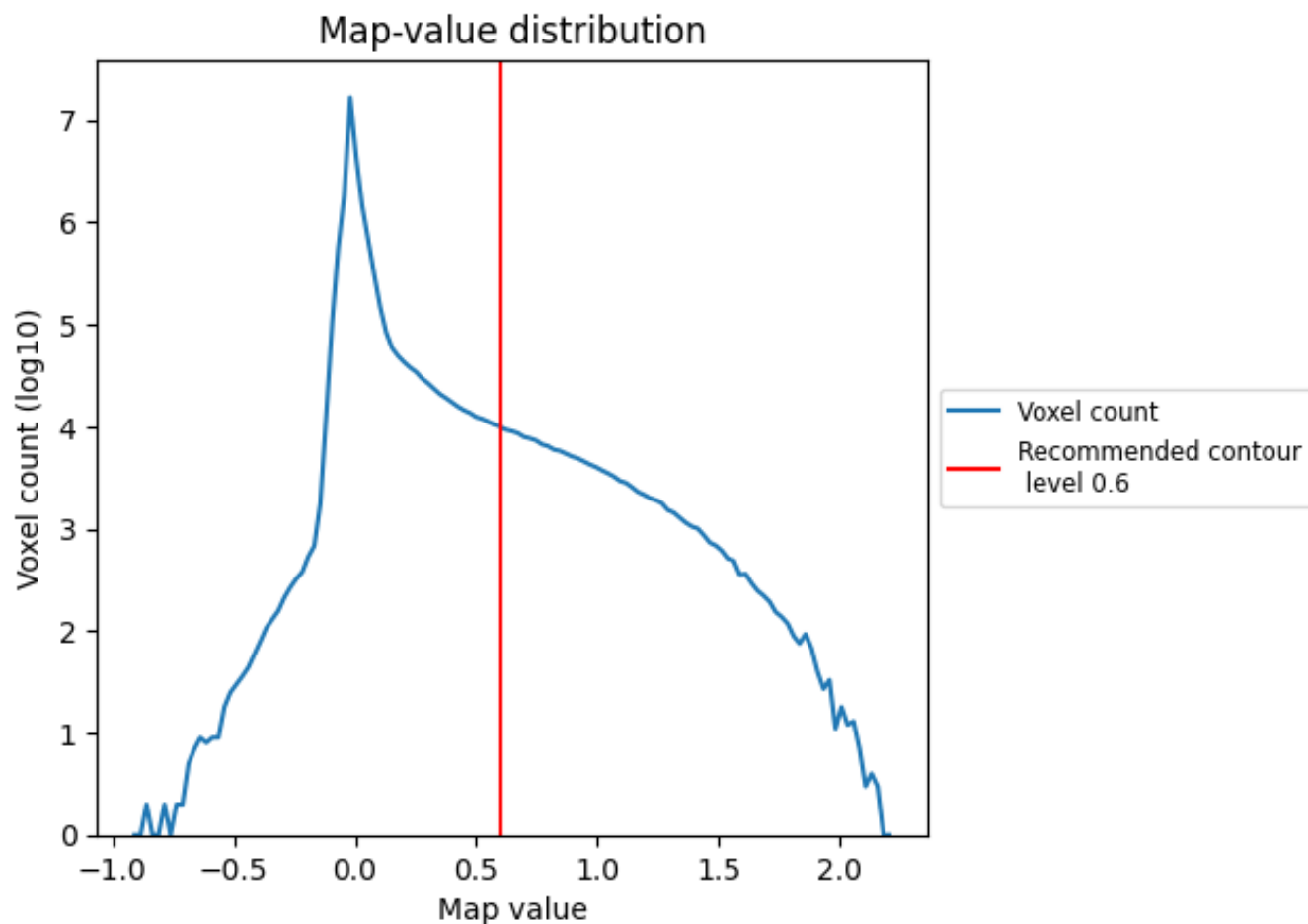


Z

## 7 Map analysis [i](#)

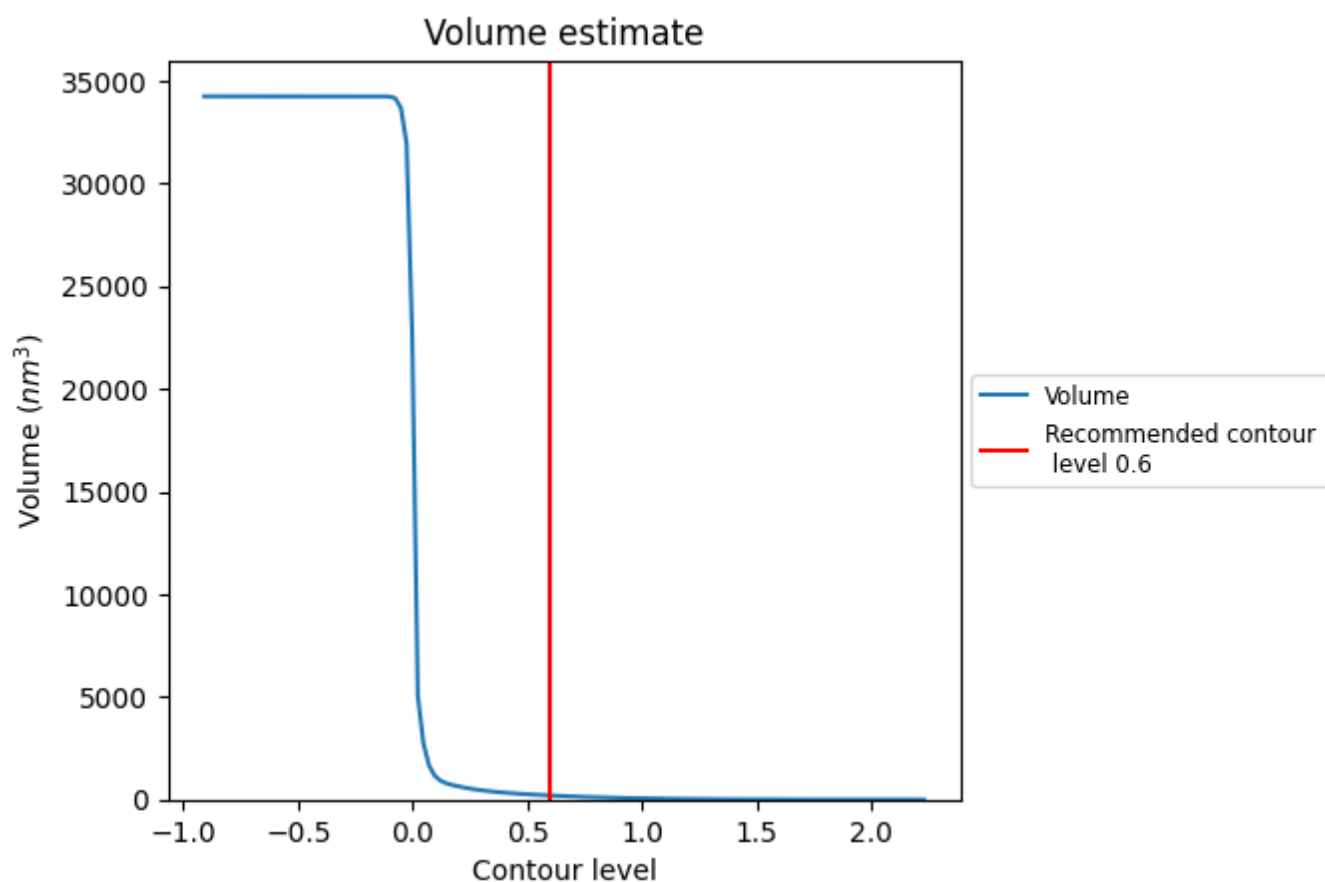
This section contains the results of statistical analysis of the map.

### 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](#)

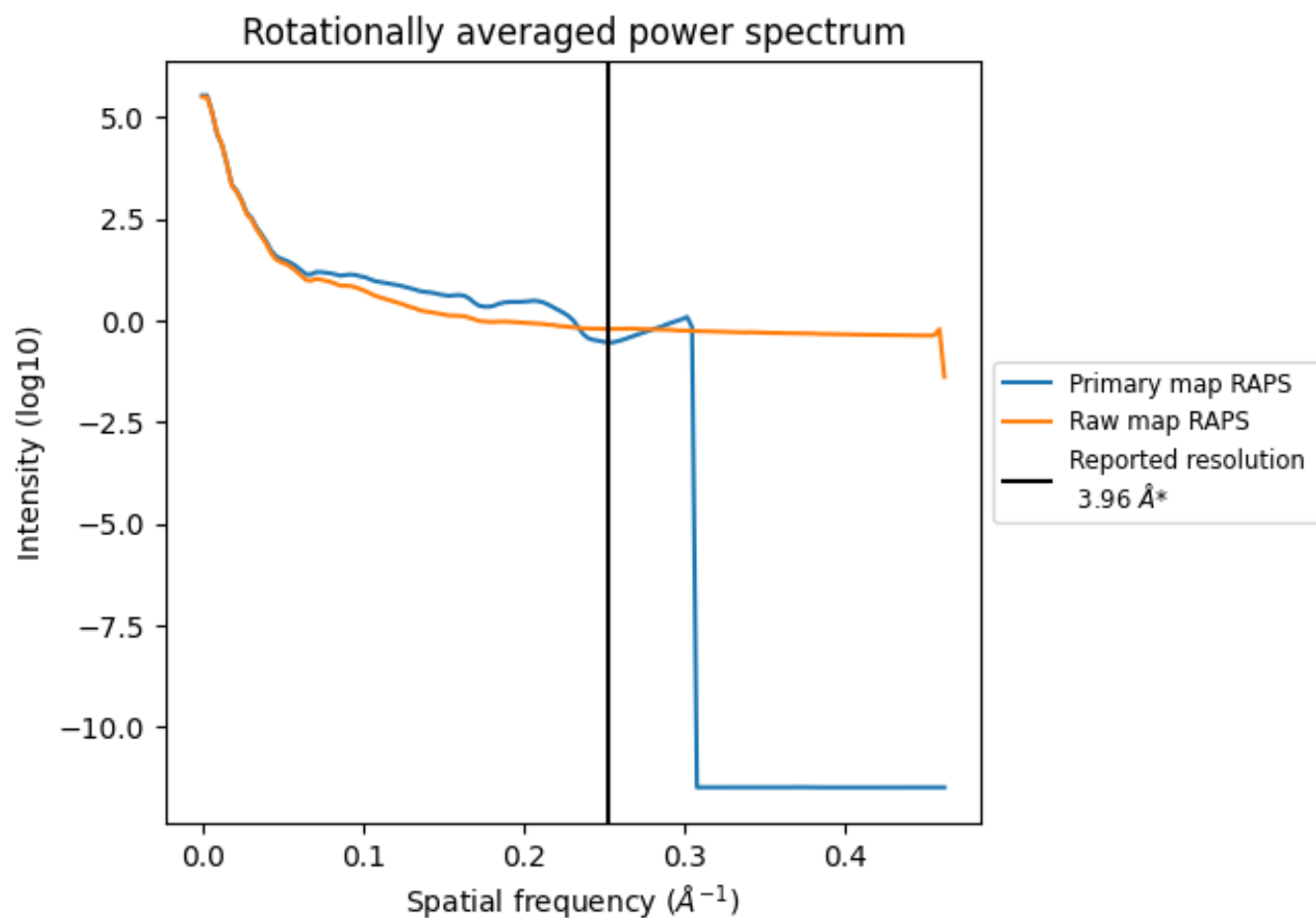


The volume at the recommended contour level is 197 nm<sup>3</sup>; this corresponds to an approximate mass of 178 kDa.

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 ⓘ

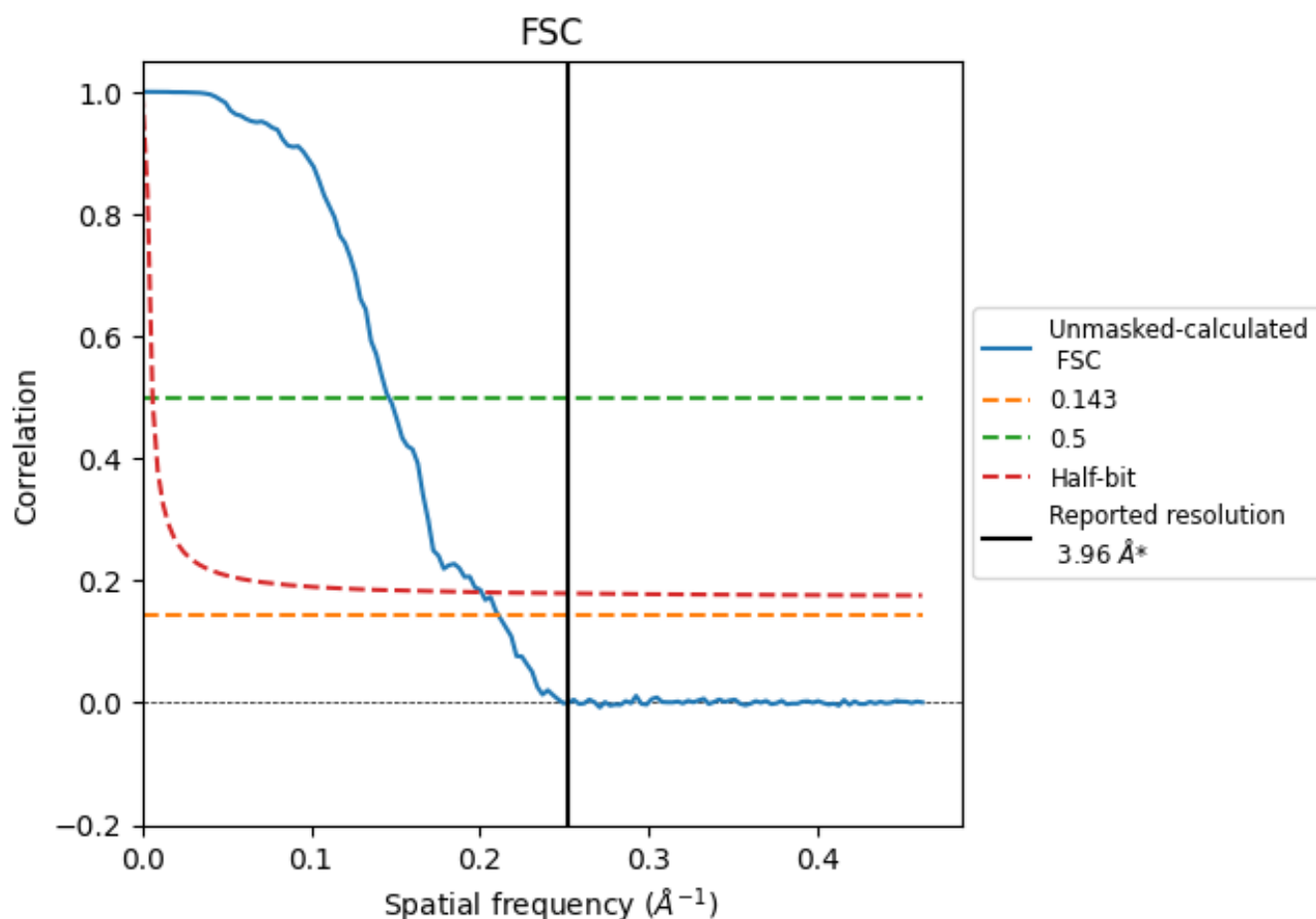


\*Reported resolution corresponds to spatial frequency of 0.253  $\text{\AA}^{-1}$

## 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 0.253  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

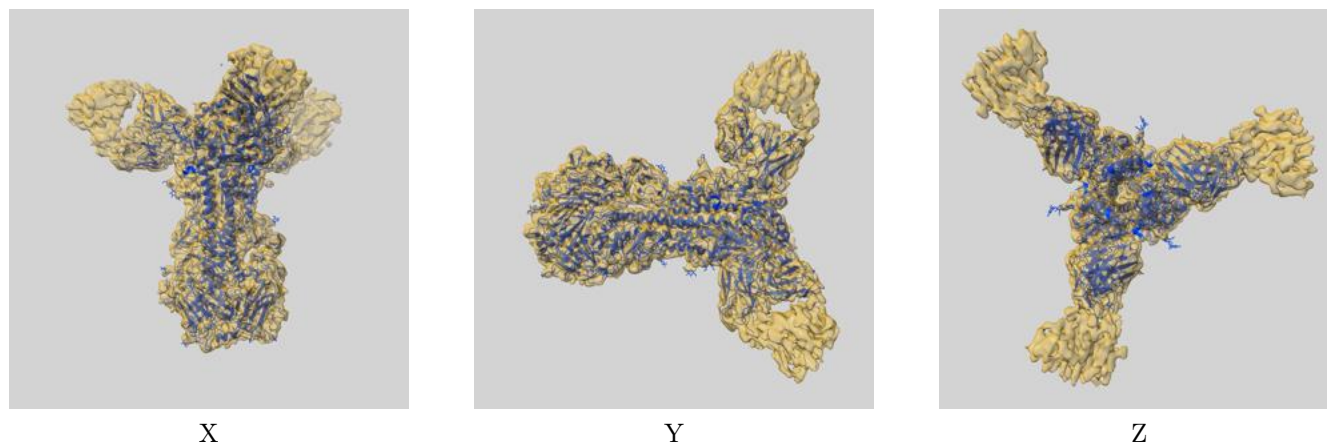
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.96	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.74	6.85	4.98

\*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 4.74 differs from the reported value 3.96 by more than 10 %

## 9 Map-model fit [i](#)

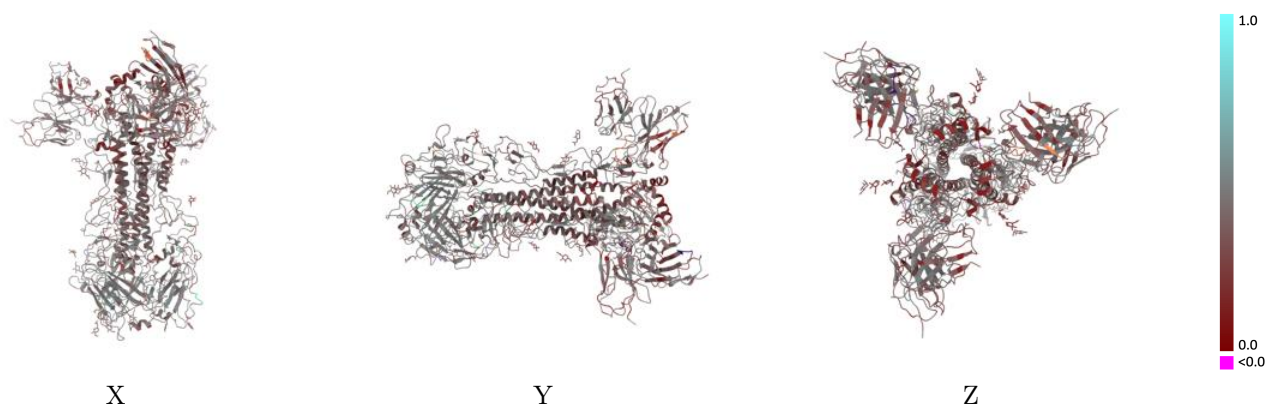
This section contains information regarding the fit between EMDB map EMD-27139 and PDB model 8D21. Per-residue inclusion information can be found in [section 3](#) on [page 11](#).

### 9.1 Map-model overlay [i](#)



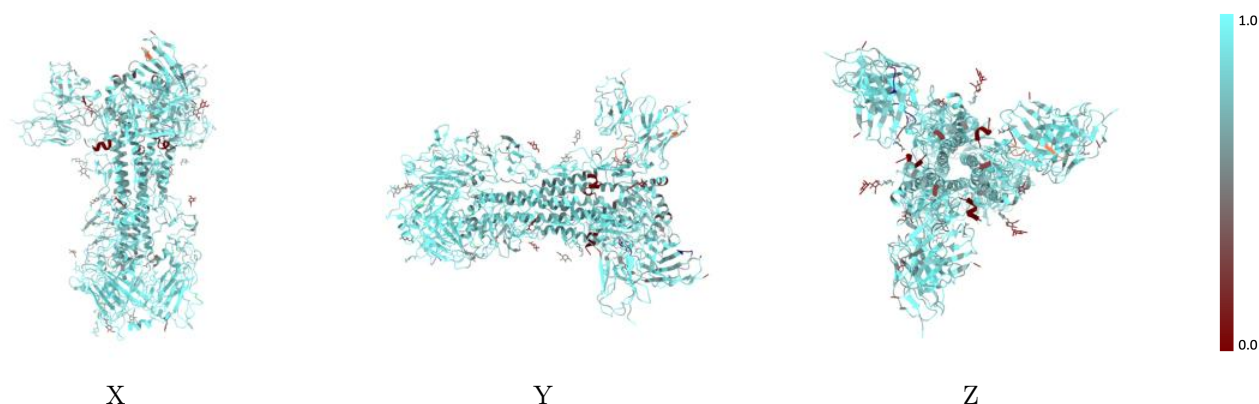
The images above show the 3D surface view of the map at the recommended contour level 0.6 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



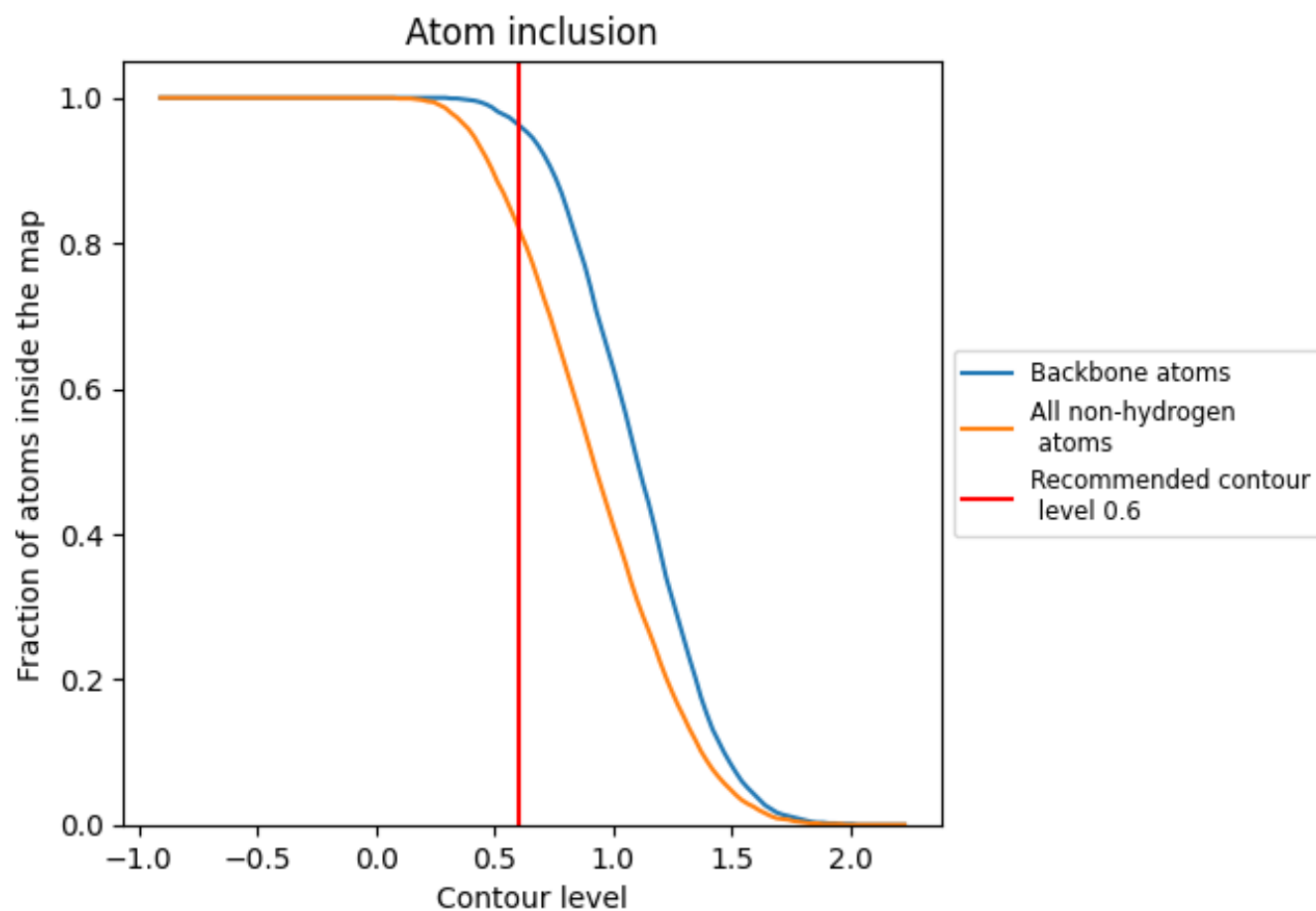
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.6).

## 9.4 Atom inclusion ⓘ



At the recommended contour level, 96% of all backbone atoms, 82% 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 (0.6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8220	<div></div> 0.3950
A	<div></div> 0.8530	<div></div> 0.4210
B	<div></div> 0.7780	<div></div> 0.3650
C	<div></div> 0.7800	<div></div> 0.3640
D	<div></div> 0.8590	<div></div> 0.4210
E	<div></div> 0.8070	<div></div> 0.3820
F	<div></div> 0.8210	<div></div> 0.3850
G	<div></div> 0.7810	<div></div> 0.3670
H	<div></div> 0.8120	<div></div> 0.3860
I	<div></div> 0.8590	<div></div> 0.4180
J	<div></div> 0.8070	<div></div> 0.3890
K	<div></div> 0.8210	<div></div> 0.3840
L	<div></div> 0.8270	<div></div> 0.3880
M	<div></div> 0.3080	<div></div> 0.2470
N	<div></div> 0.2820	<div></div> 0.2420
O	<div></div> 0.3080	<div></div> 0.2260

1.0

0.0

<0.0