



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 01:12 AM UTC

PDB ID : 8F0G / pdb_00008f0g
EMDB ID : EMD-28756
Title : Structure of SARS-CoV-2 Omicron BA.1 spike in complex with antibody Fab 1C3
Authors : Yu, X.; Zyla, D.; Hastie, K.M.; Saphire, E.O.
Deposited on : 2022-11-02
Resolution : 3.35 Å(reported)

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

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

A user guide is available at

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

The types of validation reports are described at

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

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

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

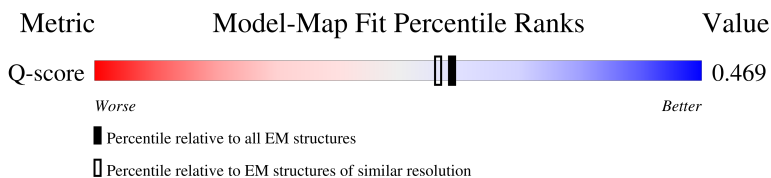
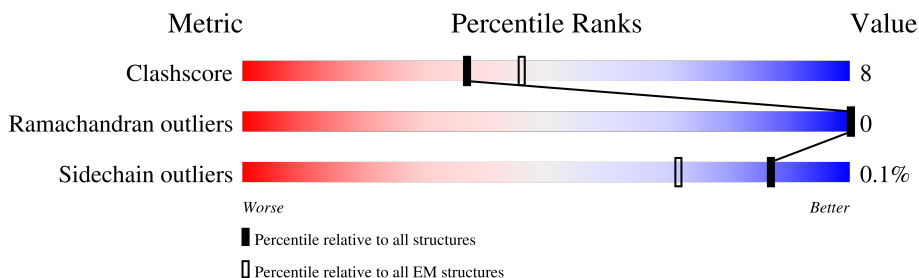
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.35 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	14390 (2.85 - 3.85)

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	X	126	9% 48% 43% 10%
2	Y	108	6% 77% 14% 9%
3	A	1209	59% 12% 29%
3	B	1209	43% 8% 49%

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Mol	Chain	Length	Quality of chain
3	C	1209	 56% 11% 34%
4	D	2	 50% 100%
4	E	2	 50% 50%
4	F	2	 50% 50%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 19635 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 Antibody 1C3 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	X	114	Total	C	N	O	S	0	0
			888	559	155	170	4		

- Molecule 2 is a protein called Antibody 1C3 Fab Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	Y	98	Total	C	N	O	S	0	0
			744	468	123	150	3		

- Molecule 3 is a protein called Spike glycoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	A	860	Total	C	N	O	S	0	0
			6684	4282	1102	1268	32		
3	B	617	Total	C	N	O	S	0	0
			4740	3005	794	917	24		
3	C	803	Total	C	N	O	S	0	0
			6243	3977	1048	1185	33		

There are 198 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	MET	-	initiating methionine	UNP P0DTC2
A	-2	GLY	-	expression tag	UNP P0DTC2
A	-1	VAL	-	expression tag	UNP P0DTC2
A	0	LYS	-	expression tag	UNP P0DTC2
A	1	VAL	-	expression tag	UNP P0DTC2
A	2	LEU	-	expression tag	UNP P0DTC2
A	3	PHE	-	expression tag	UNP P0DTC2
A	4	ALA	-	expression tag	UNP P0DTC2
A	5	LEU	-	expression tag	UNP P0DTC2
A	6	ILE	-	expression tag	UNP P0DTC2
A	7	CYS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	8	ILE	-	expression tag	UNP P0DTC2
A	9	ALA	-	expression tag	UNP P0DTC2
A	10	VAL	-	expression tag	UNP P0DTC2
A	11	ALA	-	expression tag	UNP P0DTC2
A	12	GLU	-	expression tag	UNP P0DTC2
A	13	ALA	-	expression tag	UNP P0DTC2
A	67	VAL	ALA	conflict	UNP P0DTC2
A	?	-	HIS	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	95	ILE	THR	conflict	UNP P0DTC2
A	?	-	GLY	deletion	UNP P0DTC2
A	?	-	VAL	deletion	UNP P0DTC2
A	?	-	TYR	deletion	UNP P0DTC2
A	145	ASP	TYR	conflict	UNP P0DTC2
A	211	ILE	-	insertion	UNP P0DTC2
A	211A	VAL	-	insertion	UNP P0DTC2
A	211B	ARG	ASN	conflict	UNP P0DTC2
A	211C	GLU	LEU	conflict	UNP P0DTC2
A	211D	PRO	VAL	conflict	UNP P0DTC2
A	211E	GLU	ARG	conflict	UNP P0DTC2
A	339	ASP	GLY	conflict	UNP P0DTC2
A	371	LEU	SER	conflict	UNP P0DTC2
A	373	PRO	SER	conflict	UNP P0DTC2
A	375	PHE	SER	conflict	UNP P0DTC2
A	417	ASN	LYS	conflict	UNP P0DTC2
A	440	LYS	ASN	conflict	UNP P0DTC2
A	446	SER	GLY	conflict	UNP P0DTC2
A	477	ASN	SER	conflict	UNP P0DTC2
A	478	LYS	THR	conflict	UNP P0DTC2
A	484	ALA	GLU	conflict	UNP P0DTC2
A	493	ARG	GLN	conflict	UNP P0DTC2
A	496	SER	GLY	conflict	UNP P0DTC2
A	498	ARG	GLN	conflict	UNP P0DTC2
A	501	TYR	ASN	conflict	UNP P0DTC2
A	505	HIS	TYR	conflict	UNP P0DTC2
A	547	LYS	THR	conflict	UNP P0DTC2
A	614	GLY	ASP	conflict	UNP P0DTC2
A	655	TYR	HIS	conflict	UNP P0DTC2
A	679	LYS	ASN	conflict	UNP P0DTC2
A	681	HIS	PRO	conflict	UNP P0DTC2
A	682	GLY	ARG	conflict	UNP P0DTC2
A	683	SER	ARG	conflict	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	685	SER	ARG	conflict	UNP P0DTC2
A	764	LYS	ASN	conflict	UNP P0DTC2
A	796	TYR	ASP	conflict	UNP P0DTC2
A	817	PRO	PHE	conflict	UNP P0DTC2
A	856	LYS	ASN	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	954	HIS	GLN	conflict	UNP P0DTC2
A	969	LYS	ASN	conflict	UNP P0DTC2
A	981	PHE	LEU	conflict	UNP P0DTC2
A	986	PRO	LYS	conflict	UNP P0DTC2
A	987	PRO	VAL	conflict	UNP P0DTC2
B	0	MET	-	initiating methionine	UNP P0DTC2
B	1	GLY	-	expression tag	UNP P0DTC2
B	2	VAL	-	expression tag	UNP P0DTC2
B	3	LYS	-	expression tag	UNP P0DTC2
B	4	VAL	-	expression tag	UNP P0DTC2
B	5	LEU	-	expression tag	UNP P0DTC2
B	6	PHE	-	expression tag	UNP P0DTC2
B	7	ALA	-	expression tag	UNP P0DTC2
B	8	LEU	-	expression tag	UNP P0DTC2
B	9	ILE	-	expression tag	UNP P0DTC2
B	10	CYS	-	expression tag	UNP P0DTC2
B	11	ILE	-	expression tag	UNP P0DTC2
B	12	ALA	-	expression tag	UNP P0DTC2
B	13	VAL	-	expression tag	UNP P0DTC2
B	14	ALA	-	expression tag	UNP P0DTC2
B	15	GLU	-	expression tag	UNP P0DTC2
B	16	ALA	-	expression tag	UNP P0DTC2
B	70	VAL	ALA	conflict	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	96	ILE	THR	conflict	UNP P0DTC2
B	?	-	GLY	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	?	-	TYR	deletion	UNP P0DTC2
B	143	ASP	TYR	conflict	UNP P0DTC2
B	209	ILE	-	insertion	UNP P0DTC2
B	210	VAL	-	insertion	UNP P0DTC2
B	211	ARG	ASN	conflict	UNP P0DTC2
B	212	GLU	LEU	conflict	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	213	PRO	VAL	conflict	UNP P0DTC2
B	214	GLU	ARG	conflict	UNP P0DTC2
B	339	ASP	GLY	conflict	UNP P0DTC2
B	371	LEU	SER	conflict	UNP P0DTC2
B	373	PRO	SER	conflict	UNP P0DTC2
B	375	PHE	SER	conflict	UNP P0DTC2
B	417	ASN	LYS	conflict	UNP P0DTC2
B	440	LYS	ASN	conflict	UNP P0DTC2
B	446	SER	GLY	conflict	UNP P0DTC2
B	477	ASN	SER	conflict	UNP P0DTC2
B	478	LYS	THR	conflict	UNP P0DTC2
B	484	ALA	GLU	conflict	UNP P0DTC2
B	493	ARG	GLN	conflict	UNP P0DTC2
B	496	SER	GLY	conflict	UNP P0DTC2
B	498	ARG	GLN	conflict	UNP P0DTC2
B	501	TYR	ASN	conflict	UNP P0DTC2
B	505	HIS	TYR	conflict	UNP P0DTC2
B	547	LYS	THR	conflict	UNP P0DTC2
B	614	GLY	ASP	conflict	UNP P0DTC2
B	655	TYR	HIS	conflict	UNP P0DTC2
B	679	LYS	ASN	conflict	UNP P0DTC2
B	681	HIS	PRO	conflict	UNP P0DTC2
B	682	GLY	ARG	conflict	UNP P0DTC2
B	683	SER	ARG	conflict	UNP P0DTC2
B	685	SER	ARG	conflict	UNP P0DTC2
B	764	LYS	ASN	conflict	UNP P0DTC2
B	796	TYR	ASP	conflict	UNP P0DTC2
B	817	PRO	PHE	conflict	UNP P0DTC2
B	856	LYS	ASN	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	954	HIS	GLN	conflict	UNP P0DTC2
B	969	LYS	ASN	conflict	UNP P0DTC2
B	981	PHE	LEU	conflict	UNP P0DTC2
B	986	PRO	LYS	conflict	UNP P0DTC2
B	987	PRO	VAL	conflict	UNP P0DTC2
C	0	MET	-	initiating methionine	UNP P0DTC2
C	1	GLY	-	expression tag	UNP P0DTC2
C	2	VAL	-	expression tag	UNP P0DTC2
C	3	LYS	-	expression tag	UNP P0DTC2
C	4	VAL	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	5	LEU	-	expression tag	UNP P0DTC2
C	6	PHE	-	expression tag	UNP P0DTC2
C	7	ALA	-	expression tag	UNP P0DTC2
C	8	LEU	-	expression tag	UNP P0DTC2
C	9	ILE	-	expression tag	UNP P0DTC2
C	10	CYS	-	expression tag	UNP P0DTC2
C	11	ILE	-	expression tag	UNP P0DTC2
C	12	ALA	-	expression tag	UNP P0DTC2
C	13	VAL	-	expression tag	UNP P0DTC2
C	14	ALA	-	expression tag	UNP P0DTC2
C	15	GLU	-	expression tag	UNP P0DTC2
C	16	ALA	-	expression tag	UNP P0DTC2
C	70	VAL	ALA	conflict	UNP P0DTC2
C	?	-	HIS	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	96	ILE	THR	conflict	UNP P0DTC2
C	?	-	GLY	deletion	UNP P0DTC2
C	?	-	VAL	deletion	UNP P0DTC2
C	?	-	TYR	deletion	UNP P0DTC2
C	143	ASP	TYR	conflict	UNP P0DTC2
C	209	ILE	-	insertion	UNP P0DTC2
C	210	VAL	-	insertion	UNP P0DTC2
C	211	ARG	ASN	conflict	UNP P0DTC2
C	212	GLU	LEU	conflict	UNP P0DTC2
C	213	PRO	VAL	conflict	UNP P0DTC2
C	214	GLU	ARG	conflict	UNP P0DTC2
C	339	ASP	GLY	conflict	UNP P0DTC2
C	371	LEU	SER	conflict	UNP P0DTC2
C	373	PRO	SER	conflict	UNP P0DTC2
C	375	PHE	SER	conflict	UNP P0DTC2
C	417	ASN	LYS	conflict	UNP P0DTC2
C	440	LYS	ASN	conflict	UNP P0DTC2
C	446	SER	GLY	conflict	UNP P0DTC2
C	477	ASN	SER	conflict	UNP P0DTC2
C	478	LYS	THR	conflict	UNP P0DTC2
C	484	ALA	GLU	conflict	UNP P0DTC2
C	493	ARG	GLN	conflict	UNP P0DTC2
C	496	SER	GLY	conflict	UNP P0DTC2
C	498	ARG	GLN	conflict	UNP P0DTC2
C	501	TYR	ASN	conflict	UNP P0DTC2
C	505	HIS	TYR	conflict	UNP P0DTC2
C	547	LYS	THR	conflict	UNP P0DTC2

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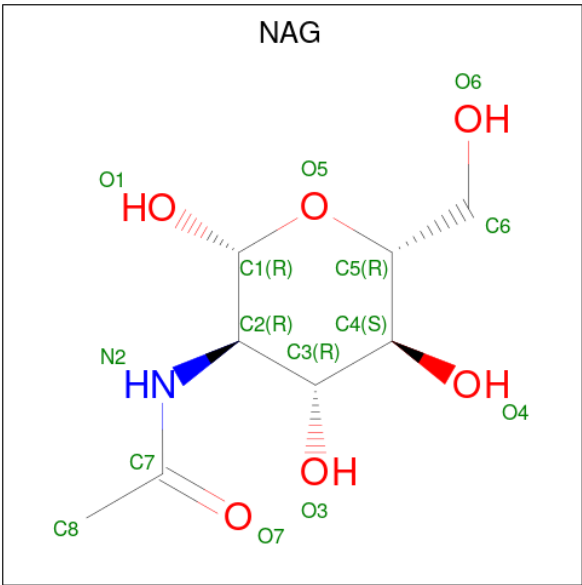
Chain	Residue	Modelled	Actual	Comment	Reference
C	614	GLY	ASP	conflict	UNP P0DTC2
C	655	TYR	HIS	conflict	UNP P0DTC2
C	679	LYS	ASN	conflict	UNP P0DTC2
C	681	HIS	PRO	conflict	UNP P0DTC2
C	682	GLY	ARG	conflict	UNP P0DTC2
C	683	SER	ARG	conflict	UNP P0DTC2
C	685	SER	ARG	conflict	UNP P0DTC2
C	764	LYS	ASN	conflict	UNP P0DTC2
C	796	TYR	ASP	conflict	UNP P0DTC2
C	817	PRO	PHE	conflict	UNP P0DTC2
C	856	LYS	ASN	conflict	UNP P0DTC2
C	892	PRO	ALA	conflict	UNP P0DTC2
C	899	PRO	ALA	conflict	UNP P0DTC2
C	942	PRO	ALA	conflict	UNP P0DTC2
C	954	HIS	GLN	conflict	UNP P0DTC2
C	969	LYS	ASN	conflict	UNP P0DTC2
C	981	PHE	LEU	conflict	UNP P0DTC2
C	986	PRO	LYS	conflict	UNP P0DTC2
C	987	PRO	VAL	conflict	UNP P0DTC2

- 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	D	2	Total	C	N	O	0	0
			28	16	2	10		
4	E	2	Total	C	N	O	0	0
			28	16	2	10		
4	F	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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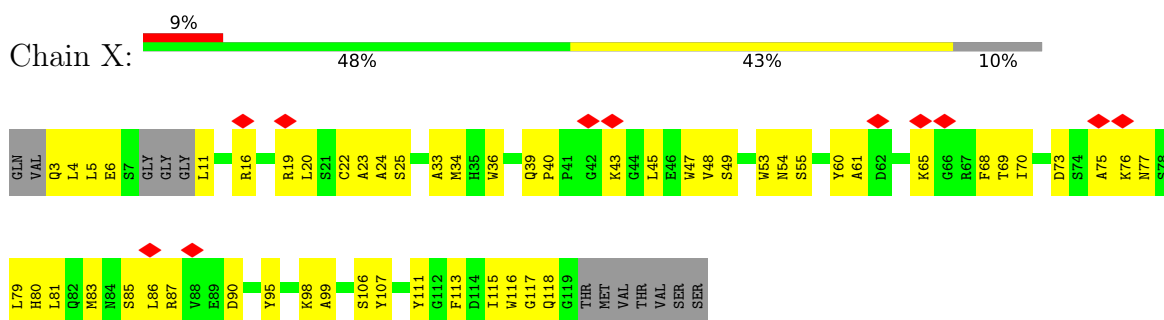
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Mol	Chain	Residues	Atoms				AltConf
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	
5	C	1	Total	C	N	O	0
			14	8	1	5	

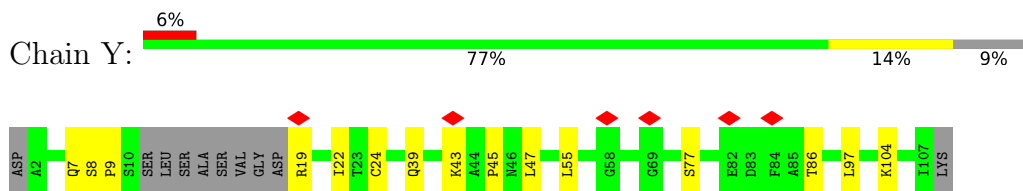
3 Residue-property plots

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

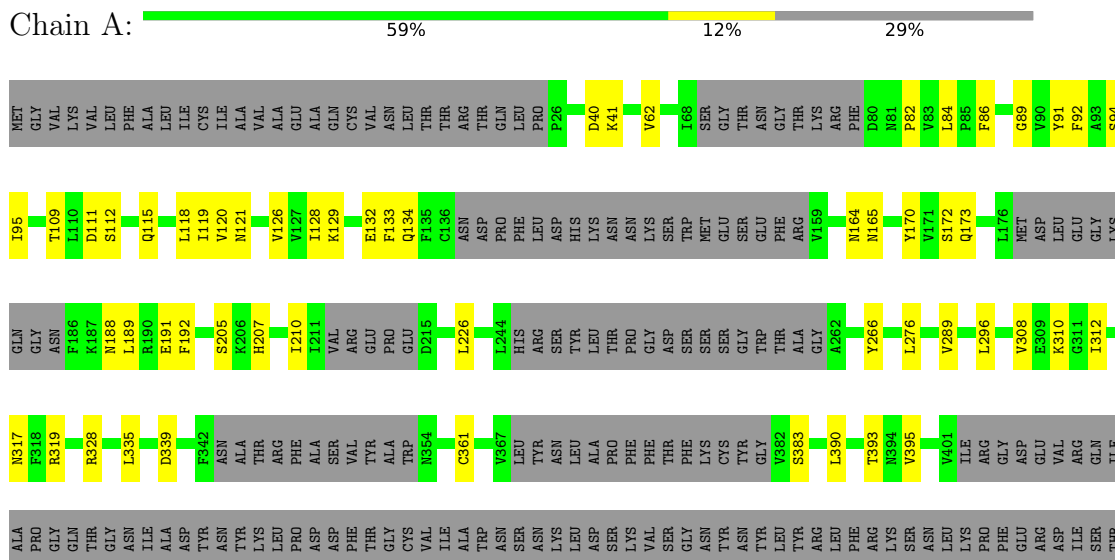
• Molecule 1: Antibody 1C3 Fab Heavy Chain



• Molecule 2: Antibody 1C3 Fab Light Chain

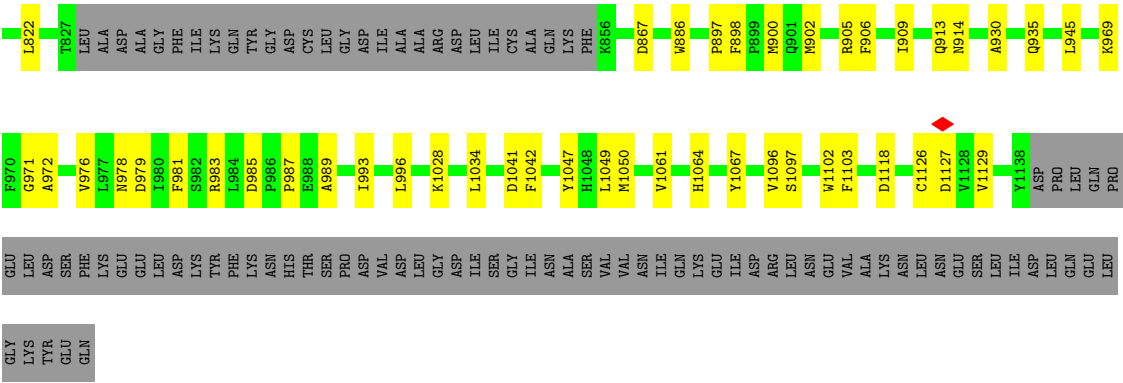


• Molecule 3: Spike glycoprotein





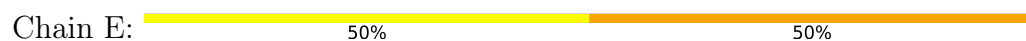
P665	ALA	PRO	LEU	VAL	THR	LYS	VAL	THR	LYS	ILE	THR	GLN	VAL	THR	GLY	PRO	MET
I666	THR	THR	PRO	ILE	ARG	LEU	PHE	LEU	GLY	ASN	LEU	GLY	VAL	THR	PHE	PHE	GLY
G667	VAL	CYS	GLU	ASP	GLY	ASP	THR	ALA	ASN	ALA	ALA	ASN	LYS	THR	ASN	SER	VAL
G669	GLY	PRO	ARG	GLU	VAL	ASP	ARG	HIS	ASN	ARG	HIS	ASN	LYS	THR	ASN	VAL	LEU
Y674	GLN	LYS	ILE	VAL	ARG	ASP	PHE	ARG	ASN	SER	THR	GLY	ASN	ASN	THR	PHE	PHE
GLN	THR	GLN	THR	ILE	GLN	SER	SER	THR	GLN	SER	THR	THR	ARG	VAL	THR	TRP	LEU
THR	GLN	S530	THR	ILE	THR	THR	VAL	THR	GLY	VAL	THR	GLY	GLU	ILE	HIS	PHE	LEU
THR	THR	LYS	ILE	PRO	ALA	THR	ALA	THR	ILE	THR	THR	THR	PHE	VAL	CYS	ILE	ILE
SER	SER	L533	TYR	GLY	ASP	ALA	SER	GLY	ASN	THR	GLY	THR	PHE	VAL	CYS	SER	ALA
HIS	THR	K537	GLN	GLN	THR	ALA	ASN	ASP	LYS	THR	ASP	ASN	LYS	GLU	THR	VAL	VAL
SER	SER	K547	GLY	GLY	THR	ALA	ASN	SER	GLY	THR	SER	SER	ILE	ASN	PHE	GLY	ALA
ALA	ALA	F559	LYS	ILE	ALA	ASN	VAL	GLY	ILE	THR	GLY	ASN	PHE	ASN	GLY	ALA	GLU
SER	SER	F563	PRO	ALA	ASP	ALA	THR	SER	THR	THR	TRP	THR	THR	ASN	LYS	CYS	GLN
VAL	VAL	Q563	CYS	ASP	ALA	CYS	ASN	THR	ASN	THR	THR	THR	PHE	ASP	ARG	VAL	VAL
ALA	ALA	D574	ASN	TYR	THR	ASN	VAL	ALA	CYS	VAL	ALA	ALA	ILE	ASN	PHE	THR	ALA
SER	SER	R577	VAL	VAL	THR	VAL	ALA	ALA	LYS	ASP	ALA	ALA	SER	ASN	THR	THR	THR
Q690	G690	R577	ALA	GLY	LEU	ALA	GLY	THR	ALA	THR	TYR	THR	LYS	ASN	VAL	ARG	ARG
I692	I692	I693	PHE	PRO	PRO	LEU	PHE	THR	ASN	THR	THR	THR	HIS	ASN	VAL	LEU	THR
I693	I693	I693	ASN	ASN	ASN	ASN	ASN	VAL	ASN	VAL	VAL	THR	LYS	ASN	PRO	GLN	GLN
G697	G697	S591	TYR	PHE	PHE	THR	THR	GLY	TYR	VAL	GLY	PRO	ILE	ASN	PHE	LEU	ASN
S698	S698	S596	PHE	THR	THR	ASN	THR	ASN	ASN	THR	THR	ASN	ILE	ASN	ASP	GLY	PRO
L699	L699	S596	PRO	GLY	GLY	LEU	GLY	LEU	THR	THR	THR	THR	ARG	THR	GLY	ALA	ALA
W703	W703	P600	LEU	ARG	VAL	ASN	VAL	ASN	CYS	ALA	Q271	VAL	VAL	THR	THR	TYR	VAL
S708	S708	Q613	SER	ILE	ILE	ASN	PHE	THR	ILE	PHE	L276	GLU	GLU	THR	PHE	ASN	THR
W709	W709	Q613	TYR	TYR	ASN	TYR	PHE	THR	ASN	THR	L276	GLU	GLU	THR	ASN	ASN	ASN
W710	W710	V622	SER	THR	THR	THR	THR	THR	THR	PHE	L285	GLU	GLU	THR	ASN	ASN	ASN
S711	S711	V622	PHE	ASN	ASN	ASN	THR	THR	THR	PHE	L285	GLU	GLU	THR	ASN	ASN	ASN
I712	I712	A633	ARG	ASN	ASN	ASN	THR	THR	THR	THR	V289	PRO	LEU	THR	ILE	ARG	ARG
W722	W722	I624	PRO	ASN	ASN	ASN	ASN	ASN	ASN	ASN	V289	GLN	GLN	THR	GLY	GLY	GLY
E725	E725	A633	THR	THR	THR	THR	THR	THR	THR	THR	D294	GLY	GLY	SER	ASN	TYR	VAL
I726	I726	ARG	GLY	GLY	GLY	GLY	ASP	VAL	ASN	VAL	D294	PHE	PHE	ASN	ILE	ILE	TYR
W740	W740	VAL	VAL	VAL	SER	VAL	SER	SER	LYS	SER	K310	ALA	ALA	ASN	ILE	PRO	ASP
Y741	Y741	THR	GLN	VAL	THR	VAL	LYS	THR	THR	THR	K310	LEU	LEU	ASN	ARG	ASP	ASP
I742	I742	GLY	THR	SER	GLY	SER	THR	THR	THR	THR	T323	CYS	CYS	GLY	GLY	LYS	VAL



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



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



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



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	49958	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)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.137	Depositor
Minimum map value	-0.342	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.020	Depositor
Recommended contour level	0.2	Depositor
Map size ($\text{\AA}$)	528.0, 528.0, 528.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.32, 1.32, 1.32	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	X	0.27	0/911	0.37	0/1234
2	Y	0.28	0/760	0.32	0/1031
3	A	0.50	0/6827	0.35	0/9288
3	B	0.49	0/4833	0.38	0/6577
3	C	0.47	0/6381	0.32	0/8678
All	All	0.48	0/19712	0.35	0/26808

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	X	888	0	839	42	0
2	Y	744	0	721	14	0
3	A	6684	0	6585	103	0
3	B	4740	0	4691	78	0
3	C	6243	0	6136	95	0
4	D	28	0	25	0	0
4	E	28	0	25	1	0
4	F	28	0	25	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	126	0	117	5	0
5	B	70	0	65	1	0
5	C	56	0	52	0	0
All	All	19635	0	19281	296	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 8.

All (296) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:641:ASN:HD22	3:C:654:GLU:HA	1.47	0.79
3:B:674:TYR:HD1	3:B:692:ILE:HG23	1.48	0.79
3:A:547:LYS:HE3	3:B:978:ASN:HB3	1.68	0.76
2:Y:9:PRO:HG2	2:Y:22:ILE:HA	1.73	0.69
1:X:16:ARG:HB2	1:X:86:LEU:HD23	1.76	0.68
3:A:62:VAL:HG11	3:A:266:TYR:HB3	1.79	0.65
3:A:886:TRP:HZ2	3:C:1107:ARG:HH12	1.42	0.65
1:X:70:ILE:HD11	1:X:79:LEU:HB3	1.78	0.65
3:A:858:LEU:HD21	3:A:962:LEU:HD23	1.78	0.65
3:B:310:LYS:HD3	3:B:600:PRO:HA	1.79	0.64
1:X:34:MET:HE2	1:X:98:LYS:HG3	1.80	0.63
3:A:577:ARG:HD3	3:A:582:LEU:HG	1.80	0.62
3:A:1102:TRP:HD1	3:A:1135:ASN:HD22	1.47	0.62
3:B:902:MET:HE3	3:B:1049:LEU:HD13	1.79	0.62
3:C:517:LEU:HD22	3:C:518:LEU:H	1.64	0.62
3:B:971:GLY:HA2	3:C:755:GLN:HE22	1.64	0.62
1:X:24:ALA:HB3	1:X:77:ASN:HB3	1.81	0.62
3:B:699:LEU:HD21	3:C:869:MET:HG2	1.82	0.61
3:A:312:ILE:HG13	3:A:598:ILE:HG13	1.82	0.61
1:X:87:ARG:HB2	1:X:90:ASP:HB2	1.81	0.61
3:A:390:LEU:HD23	3:A:390:LEU:H	1.67	0.60
3:A:707:TYR:HE1	3:B:898:PHE:H	1.47	0.60
3:B:591:SER:HB2	3:B:622:VAL:HG23	1.84	0.60
3:B:804:GLN:OE1	3:B:935:GLN:NE2	2.35	0.60
3:C:439:ASN:HA	3:C:507:PRO:HG2	1.83	0.59
3:C:398:ASP:OD1	3:C:398:ASP:N	2.32	0.59
3:A:1098:ASN:OD1	3:A:1101:HIS:N	2.31	0.59
1:X:6:GLU:HA	1:X:22:CYS:HA	1.84	0.59
1:X:40:PRO:HG2	1:X:43:LYS:HB2	1.83	0.59
1:X:47:TRP:NE1	1:X:49:SER:O	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:310:LYS:HD2	3:B:664:ILE:HD11	1.84	0.58
3:C:457:ARG:NE	3:C:467:ASP:OD2	2.33	0.58
1:X:34:MET:HE3	1:X:34:MET:N	2.18	0.58
1:X:55:SER:HB3	3:C:484:ALA:HB3	1.84	0.58
1:X:111:TYR:HB3	2:Y:97:LEU:HD11	1.85	0.58
1:X:83:MET:HA	1:X:83:MET:HE2	1.85	0.58
3:B:726:ILE:HG13	3:B:1061:VAL:HG22	1.86	0.58
1:X:48:VAL:HA	1:X:61:ALA:HB2	1.86	0.58
1:X:3:GLN:HB2	1:X:25:SER:HB3	1.86	0.58
1:X:83:MET:HB3	1:X:86:LEU:HD21	1.85	0.57
3:B:563:GLN:NE2	3:C:283:GLY:O	2.37	0.57
3:B:574:ASP:O	3:B:587:ILE:N	2.36	0.57
3:B:1047:TYR:HB2	3:B:1067:TYR:HB3	1.86	0.57
3:C:490:PHE:CE2	3:C:492:LEU:HB2	2.39	0.57
1:X:3:GLN:N	1:X:25:SER:O	2.38	0.57
3:B:742:ILE:HD11	3:B:753:LEU:HD13	1.87	0.57
3:A:308:VAL:HB	3:A:602:THR:HG23	1.86	0.57
3:A:950:ASP:O	3:A:954:HIS:ND1	2.37	0.57
5:A:1307:NAG:H2	3:B:796:TYR:CE2	2.40	0.57
3:C:897:PRO:HB2	3:C:900:MET:HG3	1.86	0.56
3:A:737:ASP:OD2	3:C:317:ASN:ND2	2.36	0.56
3:B:323:THR:HG22	3:B:324:GLU:H	1.69	0.56
3:A:627:ASP:O	3:A:634:ARG:NH2	2.37	0.56
3:A:132:GLU:H	3:A:132:GLU:CD	2.12	0.56
3:B:656:VAL:HG12	3:B:658:ASN:H	1.71	0.56
3:A:1072:GLU:HG3	5:A:1308:NAG:H62	1.86	0.55
3:B:276:LEU:HB3	3:B:289:VAL:HB	1.88	0.55
3:A:970:PHE:N	3:B:755:GLN:HE22	2.04	0.55
3:C:809:PRO:O	3:C:814:LYS:NZ	2.39	0.55
1:X:33:ALA:H	1:X:99:ALA:HB3	1.72	0.55
3:A:121:ASN:HB3	3:A:126:VAL:HG23	1.88	0.55
2:Y:7:GLN:HG2	2:Y:24:CYS:HB3	1.89	0.55
3:B:976:VAL:HG12	3:B:979:ASP:H	1.71	0.55
2:Y:39:GLN:HA	2:Y:45:PRO:HA	1.89	0.55
1:X:4:LEU:H	1:X:118:GLN:HE22	1.55	0.54
1:X:95:TYR:HE2	2:Y:45:PRO:HD2	1.71	0.54
3:A:599:THR:HG22	3:A:608:VAL:HG12	1.88	0.54
3:B:547:LYS:HB2	3:C:978:ASN:HB3	1.89	0.54
3:A:708:SER:O	3:A:711:SER:HB3	2.07	0.54
3:A:1076:THR:OG1	3:A:1097:SER:OG	2.25	0.54
3:C:1126:CYS:SG	3:C:1127:ASP:N	2.81	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:39:GLN:HB2	1:X:45:LEU:HD23	1.89	0.54
3:B:989:ALA:O	3:B:993:ILE:HD12	2.07	0.54
3:A:726:ILE:HG12	3:A:1061:VAL:HG22	1.90	0.54
3:A:985:ASP:HB3	3:A:987:PRO:HD2	1.90	0.54
1:X:85:SER:O	1:X:87:ARG:NH1	2.40	0.54
3:A:964:LYS:NZ	3:C:569:ILE:O	2.40	0.53
3:B:669:GLY:HA3	3:C:869:MET:HE1	1.88	0.53
3:A:1086:LYS:HE2	3:A:1122:VAL:HG11	1.89	0.53
1:X:70:ILE:HD13	1:X:81:LEU:HD12	1.90	0.53
3:A:112:SER:HA	3:A:132:GLU:HB2	1.91	0.53
2:Y:86:THR:HG22	2:Y:104:LYS:HB3	1.91	0.53
3:B:662:CYS:HB2	3:B:697:MET:HE3	1.91	0.53
1:X:19:ARG:HH21	1:X:80:HIS:HB3	1.74	0.53
3:C:675:GLN:O	3:C:690:GLN:N	2.42	0.53
3:A:188:ASN:OD1	3:A:207:HIS:NE2	2.41	0.52
3:A:115:GLN:HA	3:A:132:GLU:HA	1.90	0.52
1:X:116:TRP:CD2	2:Y:45:PRO:HG2	2.45	0.52
3:A:310:LYS:HG3	3:A:600:PRO:HA	1.92	0.52
3:A:722:VAL:HG22	3:A:930:ALA:HB1	1.92	0.52
3:A:811:LYS:HG2	3:A:813:SER:H	1.74	0.52
3:B:646:ARG:HB2	3:B:668:ALA:HB1	1.92	0.52
1:X:36:TRP:HB2	1:X:49:SER:HB2	1.91	0.52
3:B:559:PHE:HD2	3:B:577:ARG:HH21	1.56	0.52
3:C:498:ARG:HD3	3:C:499:PRO:HD2	1.91	0.52
3:A:1123:SER:HB3	3:B:914:ASN:HD22	1.75	0.51
3:B:971:GLY:CA	3:C:755:GLN:HE22	2.23	0.51
3:C:1106:GLN:NE2	3:C:1111:GLU:OE1	2.43	0.51
3:C:709:ASN:OD1	3:C:709:ASN:N	2.42	0.51
1:X:20:LEU:O	1:X:81:LEU:N	2.43	0.51
3:A:1090:PRO:HD3	3:A:1095:PHE:CE2	2.46	0.51
1:X:73:ASP:OD2	1:X:80:HIS:NE2	2.41	0.51
3:B:654:GLU:HG3	3:B:693:ILE:HG22	1.92	0.51
3:C:1028:LYS:NZ	3:C:1042:PHE:O	2.44	0.51
3:A:1029:MET:HE1	3:A:1053:PRO:HB3	1.91	0.51
3:B:708:SER:O	3:B:711:SER:OG	2.24	0.51
1:X:5:LEU:O	1:X:23:ALA:N	2.43	0.51
3:C:985:ASP:OD1	3:C:985:ASP:N	2.44	0.51
3:A:1125:ASN:ND2	3:A:1127:ASP:OD2	2.44	0.51
3:A:762:GLN:N	3:A:762:GLN:OE1	2.44	0.50
1:X:115:ILE:HG22	1:X:117:GLY:H	1.76	0.50
3:C:738:CYS:SG	3:C:739:THR:N	2.84	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:75:ALA:HB3	1:X:76:LYS:HD3	1.92	0.50
3:A:1135:ASN:OD1	3:A:1136:THR:N	2.45	0.50
3:B:285:ILE:HD12	3:B:285:ILE:O	2.11	0.50
3:C:277:LEU:HB3	3:C:285:ILE:HD12	1.93	0.50
1:X:54:ASN:HB3	3:C:484:ALA:HB1	1.94	0.50
3:A:890:ALA:HB1	3:C:1046:GLY:HA2	1.94	0.50
3:C:708:SER:HB2	3:C:711:SER:HB3	1.94	0.50
3:A:762:GLN:NE2	3:C:961:THR:OG1	2.45	0.50
3:A:1029:MET:HA	3:A:1029:MET:HE3	1.93	0.49
3:C:1118:ASP:OD1	3:C:1119:ASN:N	2.44	0.49
1:X:11:LEU:N	1:X:16:ARG:HE	2.10	0.49
3:B:906:PHE:O	3:B:909:ILE:HG12	2.13	0.49
3:C:642:VAL:HG12	3:C:651:ILE:HD11	1.93	0.49
3:A:393:THR:HB	3:A:522:ALA:HA	1.95	0.49
3:B:674:TYR:HA	3:B:692:ILE:HA	1.93	0.48
1:X:4:LEU:HB2	1:X:117:GLY:HA2	1.95	0.48
3:A:808:ASP:OD1	3:A:808:ASP:N	2.44	0.48
3:C:461:LEU:HD12	3:C:465:GLU:HB2	1.95	0.48
3:C:374:PHE:H	3:C:436:TRP:HB3	1.77	0.48
1:X:106:SER:OG	1:X:107:TYR:N	2.46	0.48
3:A:1089:PHE:HB3	3:B:913:GLN:NE2	2.29	0.48
3:B:710:ASN:ND2	3:B:710:ASN:O	2.47	0.48
3:B:722:VAL:HG22	3:B:930:ALA:HB1	1.94	0.48
3:A:118:LEU:HD11	3:A:120:VAL:HB	1.94	0.48
3:C:851:CYS:SG	3:C:852:ALA:N	2.85	0.48
3:A:82:PRO:HG2	3:A:84:LEU:HD21	1.96	0.48
3:A:733:LYS:HE3	3:A:861:LEU:HB2	1.96	0.48
2:Y:39:GLN:HB2	2:Y:45:PRO:HB3	1.95	0.47
3:B:969:LYS:HA	3:B:969:LYS:HE2	1.96	0.47
3:B:978:ASN:HA	3:B:981:PHE:CD2	2.50	0.47
3:A:94:SER:C	3:A:95:ILE:HD13	2.39	0.47
3:A:1089:PHE:HB3	3:B:913:GLN:HE21	1.80	0.47
3:C:460:ASN:OD1	3:C:460:ASN:N	2.47	0.47
3:A:109:THR:OG1	3:A:111:ASP:OD1	2.28	0.47
3:A:319:ARG:NH2	3:A:592:PHE:O	2.42	0.47
3:A:132:GLU:OE2	3:A:165:ASN:HB2	2.14	0.47
3:A:335:LEU:O	3:A:361:CYS:HB2	2.14	0.47
3:B:712:ILE:HD13	3:C:896:ILE:HD13	1.96	0.47
3:B:1129:VAL:HG13	3:C:917:TYR:HB3	1.97	0.47
3:C:518:LEU:C	3:C:519:HIS:HD1	2.23	0.47
3:A:112:SER:HB3	3:A:134:GLN:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:666:ILE:HD11	3:C:672:ALA:HB2	1.97	0.47
3:C:986:PRO:N	3:C:987:PRO:HD2	2.30	0.47
1:X:68:PHE:HD1	1:X:83:MET:HE1	1.80	0.47
3:C:578:ASP:OD1	3:C:578:ASP:N	2.47	0.47
3:C:409:GLN:HB3	3:C:419:ALA:HB2	1.98	0.46
3:C:316:SER:OG	3:C:317:ASN:N	2.47	0.46
3:C:490:PHE:HE2	3:C:492:LEU:HB2	1.77	0.46
3:B:822:LEU:HD22	3:B:945:LEU:HD21	1.98	0.46
3:B:1096:VAL:O	3:B:1103:PHE:N	2.42	0.46
2:Y:43:LYS:HB3	2:Y:43:LYS:HE3	1.73	0.46
3:A:132:GLU:OE1	3:A:164:ASN:N	2.48	0.46
3:B:867:ASP:OD1	3:B:867:ASP:N	2.49	0.46
3:A:900:MET:HE2	3:A:900:MET:HB3	1.77	0.46
3:C:1108:ASN:OD1	3:C:1108:ASN:N	2.48	0.46
3:A:119:ILE:HG23	3:A:128:ILE:HG12	1.98	0.45
3:A:317:ASN:OD1	3:A:317:ASN:N	2.47	0.45
3:C:885:GLY:O	3:C:887:THR:HG23	2.16	0.45
3:C:1029:MET:HE1	3:C:1053:PRO:HB3	1.98	0.45
3:A:1083:HIS:CD2	3:A:1137:VAL:HB	2.51	0.45
3:B:725:GLU:OE1	3:B:1064:HIS:NE2	2.49	0.45
3:A:62:VAL:CG1	3:A:266:TYR:HB3	2.45	0.45
3:A:707:TYR:HE1	3:B:898:PHE:N	2.14	0.45
3:A:612:TYR:O	3:A:648:GLY:HA3	2.16	0.45
3:A:129:LYS:HB3	3:A:133:PHE:HZ	1.82	0.45
3:C:752:LEU:O	3:C:756:TYR:HD1	2.00	0.45
3:A:328:ARG:HD3	3:A:543:PHE:CE2	2.51	0.45
3:C:290:ASP:O	3:C:297:SER:HB3	2.17	0.45
3:A:1100:THR:OG1	3:A:1101:HIS:ND1	2.50	0.44
3:A:86:PHE:CZ	3:A:89:GLY:HA2	2.52	0.44
3:B:591:SER:HB2	3:B:622:VAL:CG2	2.46	0.44
3:B:985:ASP:HB3	3:B:987:PRO:HD2	1.99	0.44
1:X:45:LEU:HD13	1:X:116:TRP:CH2	2.52	0.44
3:B:755:GLN:O	3:B:755:GLN:NE2	2.50	0.44
3:A:606:ASN:O	3:A:608:VAL:HG13	2.15	0.44
1:X:68:PHE:HA	1:X:83:MET:HE1	1.99	0.44
3:C:331:ASN:OD1	3:C:331:ASN:N	2.50	0.44
3:C:735:SER:O	3:C:859:THR:OG1	2.31	0.44
3:A:973:ILE:HD12	3:A:984:LEU:HD11	2.00	0.44
3:A:985:ASP:N	3:C:383:SER:OG	2.50	0.44
3:C:470:THR:HB	3:C:490:PHE:CE1	2.53	0.44
3:C:518:LEU:O	3:C:519:HIS:ND1	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:733:LYS:NZ	3:C:775:ASP:OD2	2.39	0.44
3:C:767:LEU:HD23	3:C:767:LEU:HA	1.82	0.44
2:Y:19:ARG:HD2	2:Y:77:SER:HA	1.99	0.44
3:A:1106:GLN:HG3	3:A:1109:PHE:O	2.17	0.44
3:B:981:PHE:HE1	3:B:993:ILE:HG12	1.83	0.44
3:C:347:PHE:HB2	3:C:401:VAL:HG23	1.99	0.43
3:C:661:GLU:O	3:C:695:TYR:OH	2.26	0.43
3:B:1097:SER:HB2	3:B:1102:TRP:CD2	2.54	0.43
3:C:366:SER:HA	3:C:369:TYR:HD2	1.82	0.43
3:B:537:LYS:NZ	3:B:624:ILE:HG21	2.32	0.43
3:B:709:ASN:HA	3:C:897:PRO:HG3	1.99	0.43
3:C:310:LYS:HG3	3:C:600:PRO:HA	2.00	0.43
3:C:900:MET:HE2	3:C:900:MET:HB3	1.73	0.43
3:A:970:PHE:H	3:B:755:GLN:HE22	1.63	0.43
3:B:969:LYS:NZ	3:B:972:ALA:O	2.49	0.43
3:C:406:GLU:OE1	3:C:495:TYR:OH	2.32	0.43
1:X:116:TRP:CG	2:Y:45:PRO:HG2	2.54	0.43
3:A:92:PHE:O	3:A:192:PHE:N	2.48	0.43
3:A:773:GLU:OE2	3:A:1019:ARG:NH1	2.51	0.43
3:A:1098:ASN:HB2	5:A:1309:NAG:H2	2.00	0.43
3:B:294:ASP:HB2	3:B:295:PRO:HD2	2.00	0.43
3:B:996:LEU:HD23	3:B:996:LEU:HA	1.80	0.43
3:B:1028:LYS:NZ	3:B:1042:PHE:O	2.51	0.43
3:A:128:ILE:HD12	3:A:170:TYR:HD2	1.84	0.43
3:A:1000:ARG:O	3:A:1003:SER:OG	2.36	0.43
3:C:346:ARG:HD2	3:C:346:ARG:HA	1.94	0.43
3:C:651:ILE:HD13	3:C:651:ILE:HA	1.88	0.43
3:C:726:ILE:HG12	3:C:1061:VAL:HG22	2.01	0.43
3:C:819:GLU:OE2	3:C:1055:SER:OG	2.30	0.43
3:C:368:LEU:HD12	3:C:368:LEU:O	2.18	0.43
3:C:478:LYS:HD2	3:C:479:PRO:HD2	2.01	0.43
3:B:905:ARG:NE	3:B:1050:MET:HE3	2.34	0.42
3:C:517:LEU:HD13	3:C:518:LEU:N	2.34	0.42
3:C:329:PHE:O	3:C:580:GLN:NE2	2.51	0.42
3:C:1097:SER:HB2	3:C:1102:TRP:CD2	2.55	0.42
2:Y:55:LEU:HD13	2:Y:55:LEU:HA	1.93	0.42
3:B:1049:LEU:HD23	3:B:1049:LEU:HA	1.84	0.42
3:C:616:ASN:OD1	3:C:618:THR:N	2.37	0.42
3:C:964:LYS:HD2	3:C:964:LYS:HA	1.66	0.42
1:X:60:TYR:HB2	1:X:65:LYS:NZ	2.34	0.42
3:A:906:PHE:HE1	3:A:1049:LEU:HD11	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:703:ASN:ND2	3:C:787:GLN:OE1	2.53	0.42
3:B:1041:ASP:HB3	3:C:1030:SER:HB3	2.00	0.42
3:C:806:LEU:HD23	3:C:806:LEU:HA	1.86	0.42
3:A:905:ARG:NH1	3:A:1049:LEU:O	2.51	0.42
3:C:804:GLN:NE2	4:F:1:NAG:O6	2.38	0.42
3:A:205:SER:HB3	3:A:226:LEU:HD22	2.01	0.42
3:A:296:LEU:HD13	3:A:608:VAL:HG11	2.02	0.42
3:C:804:GLN:HE21	4:F:1:NAG:HO6	1.61	0.42
3:C:884:SER:OG	3:C:887:THR:OG1	2.38	0.42
1:X:113:PHE:O	2:Y:47:LEU:HD11	2.19	0.42
3:A:91:TYR:OH	3:A:191:GLU:OE2	2.32	0.42
3:A:675:GLN:NE2	3:A:690:GLN:HB2	2.35	0.42
3:B:1126:CYS:SG	3:B:1127:ASP:N	2.93	0.42
3:C:529:LYS:HD2	3:C:529:LYS:HA	1.88	0.42
3:C:722:VAL:HG22	3:C:1065:VAL:HG22	2.02	0.42
4:E:1:NAG:H62	4:E:2:NAG:H83	2.01	0.42
3:A:41:LYS:HD2	3:C:520:ALA:HB2	2.01	0.41
3:A:764:LYS:HB2	3:A:764:LYS:HE3	1.76	0.41
3:A:1004:LEU:HD23	3:A:1004:LEU:HA	1.88	0.41
3:A:395:VAL:HG22	3:A:515:PHE:HB3	2.02	0.41
3:A:934:ILE:HA	3:A:934:ILE:HD13	1.84	0.41
3:B:740:MET:HE3	3:B:740:MET:HB2	1.83	0.41
3:B:1103:PHE:CE2	5:B:1304:NAG:H62	2.55	0.41
3:A:189:LEU:HB2	3:A:210:ILE:HD11	2.03	0.41
3:A:276:LEU:HB3	3:A:289:VAL:HG22	2.02	0.41
3:A:1001:LEU:HD23	3:A:1001:LEU:HA	1.91	0.41
3:A:1052:PHE:HB2	3:A:1063:LEU:HB2	2.02	0.41
1:X:60:TYR:CZ	1:X:69:THR:HA	2.54	0.41
3:A:172:SER:OG	3:A:173:GLN:N	2.51	0.41
3:A:600:PRO:HB3	3:A:674:TYR:HB2	2.03	0.41
3:C:750:SER:O	3:C:753:LEU:HG	2.21	0.41
3:A:966:LEU:HD23	3:A:966:LEU:HA	1.87	0.41
3:C:353:TRP:CD1	3:C:353:TRP:H	2.39	0.41
3:C:596:SER:O	3:C:611:LEU:N	2.49	0.41
3:C:752:LEU:HG	3:C:993:ILE:HD11	2.02	0.41
1:X:53:TRP:CD1	1:X:53:TRP:H	2.39	0.41
3:B:886:TRP:HB2	3:B:1034:LEU:O	2.21	0.41
3:B:897:PRO:HG2	3:B:900:MET:HG3	2.03	0.41
3:C:296:LEU:HD23	3:C:608:VAL:HG11	2.02	0.41
3:B:533:LEU:HD23	3:B:533:LEU:H	1.86	0.41
3:A:629:LEU:HD23	3:A:629:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:964:LYS:NZ	3:C:569:ILE:HG13	2.35	0.41
5:A:1307:NAG:H61	3:B:796:TYR:HB3	2.02	0.41
3:B:596:SER:OG	3:B:613:GLN:NE2	2.54	0.41
3:C:410:ILE:HD13	3:C:510:VAL:HG11	2.03	0.41
3:A:984:LEU:HB2	3:A:989:ALA:HB2	2.03	0.40
3:B:1118:ASP:N	3:B:1118:ASP:OD1	2.52	0.40
3:C:425:LEU:HD23	3:C:425:LEU:HA	1.88	0.40
2:Y:8:SER:HB3	2:Y:9:PRO:HD3	2.02	0.40
3:A:40:ASP:OD1	3:A:40:ASP:N	2.53	0.40
3:A:339:ASP:OD1	3:A:339:ASP:N	2.48	0.40
3:A:703:ASN:O	3:B:789:TYR:HA	2.21	0.40
3:A:1090:PRO:HD2	3:B:913:GLN:HE22	1.86	0.40
3:B:979:ASP:OD1	3:B:983:ARG:HG3	2.20	0.40
3:A:383:SER:OG	3:B:985:ASP:OD1	2.30	0.40
3:A:390:LEU:H	3:A:390:LEU:CD2	2.33	0.40
3:C:878:LEU:HA	3:C:878:LEU:HD23	1.85	0.40
3:A:909:ILE:HG13	3:A:911:VAL:HG23	2.03	0.40
5:A:1307:NAG:H2	3:B:796:TYR:HE2	1.84	0.40
3:B:802:PHE:HD1	3:B:805:ILE:HD11	1.87	0.40
3:C:752:LEU:HG	3:C:993:ILE:CD1	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	X	110/126 (87%)	103 (94%)	7 (6%)	0	100	100
2	Y	94/108 (87%)	85 (90%)	9 (10%)	0	100	100
3	A	834/1209 (69%)	802 (96%)	32 (4%)	0	100	100
3	B	607/1209 (50%)	585 (96%)	22 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	793/1209 (66%)	776 (98%)	17 (2%)	0	100	100
All	All	2438/3861 (63%)	2351 (96%)	87 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	X	95/104 (91%)	95 (100%)	0	100	100
2	Y	84/92 (91%)	84 (100%)	0	100	100
3	A	756/1055 (72%)	756 (100%)	0	100	100
3	B	535/1055 (51%)	534 (100%)	1 (0%)	87	85
3	C	698/1055 (66%)	696 (100%)	2 (0%)	86	84
All	All	2168/3361 (64%)	2165 (100%)	3 (0%)	87	88

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	B	666	ILE
3	C	488	CYS
3	C	751	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	39	GLN
2	Y	46	ASN
2	Y	80	GLN
3	A	690	GLN
3	A	755	GLN
3	A	774	GLN

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Mol	Chain	Res	Type
3	A	787	GLN
3	A	1083	HIS
3	B	280	ASN
3	B	282	ASN
3	B	563	GLN
3	B	613	GLN
3	B	755	GLN
3	B	913	GLN
3	B	1010	GLN
3	C	542	ASN
3	C	641	ASN
3	C	926	GLN

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 ⓘ

6 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	3,4	14,14,15	0.21	0	17,19,21	0.43	0
4	NAG	D	2	4	14,14,15	0.18	0	17,19,21	0.44	0
4	NAG	E	1	3,4	14,14,15	0.65	1 (7%)	17,19,21	0.53	0
4	NAG	E	2	4	14,14,15	0.47	0	17,19,21	0.57	0
4	NAG	F	1	3,4	14,14,15	0.25	0	17,19,21	0.50	0
4	NAG	F	2	4	14,14,15	0.31	0	17,19,21	0.56	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	3,4	-	2/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1
4	NAG	E	1	3,4	-	3/6/23/26	0/1/1/1
4	NAG	E	2	4	-	4/6/23/26	0/1/1/1
4	NAG	F	1	3,4	-	2/6/23/26	0/1/1/1
4	NAG	F	2	4	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1	NAG	O5-C1	-2.21	1.40	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (15) torsion outliers are listed below:

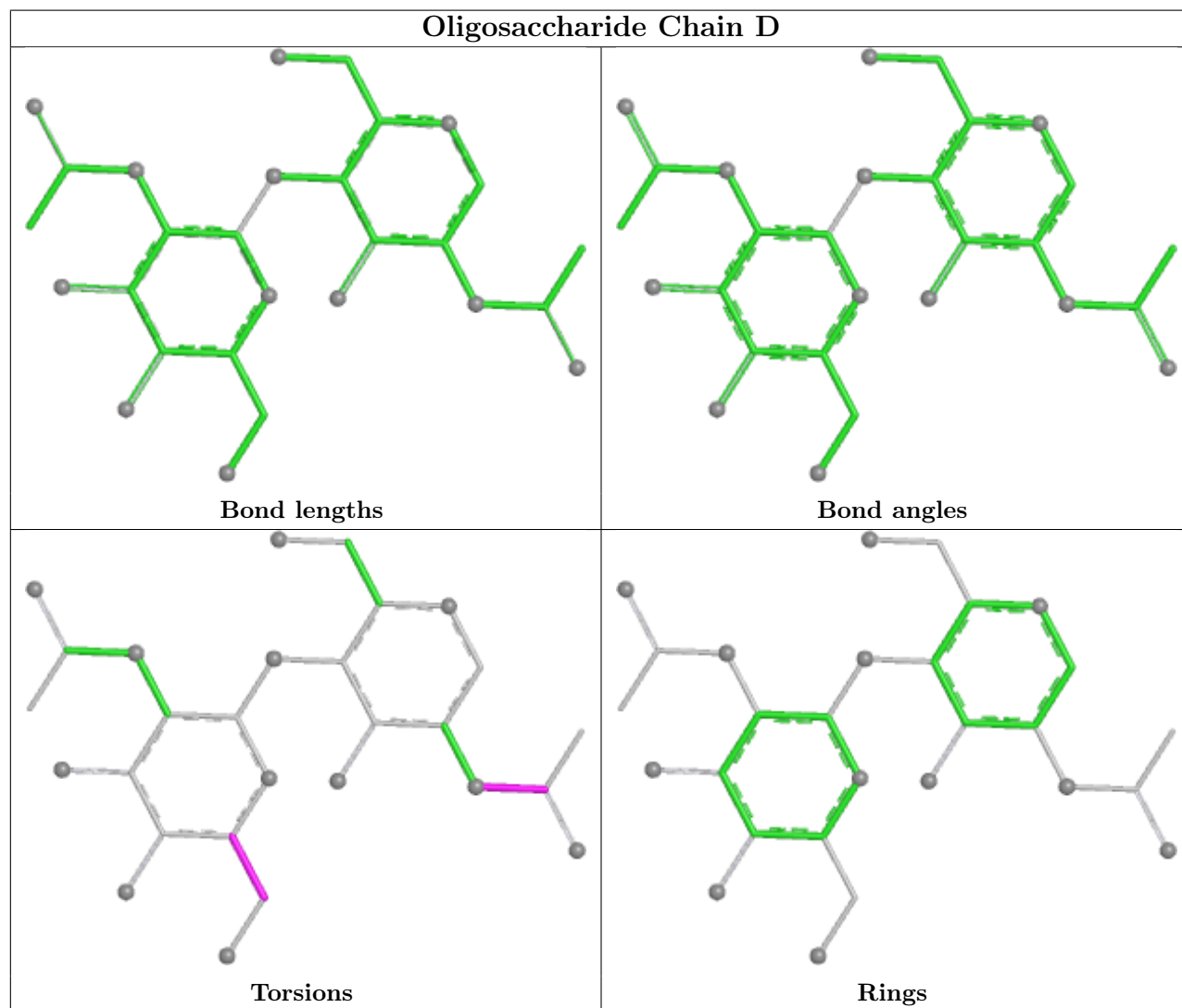
Mol	Chain	Res	Type	Atoms
4	D	2	NAG	O5-C5-C6-O6
4	D	2	NAG	C4-C5-C6-O6
4	D	1	NAG	C8-C7-N2-C2
4	D	1	NAG	O7-C7-N2-C2
4	F	1	NAG	C8-C7-N2-C2
4	F	1	NAG	O7-C7-N2-C2
4	E	2	NAG	O5-C5-C6-O6
4	E	1	NAG	O5-C5-C6-O6
4	E	1	NAG	C1-C2-N2-C7
4	E	2	NAG	C3-C2-N2-C7
4	F	2	NAG	C3-C2-N2-C7
4	E	2	NAG	C1-C2-N2-C7
4	F	2	NAG	C1-C2-N2-C7
4	E	1	NAG	C3-C2-N2-C7
4	E	2	NAG	C4-C5-C6-O6

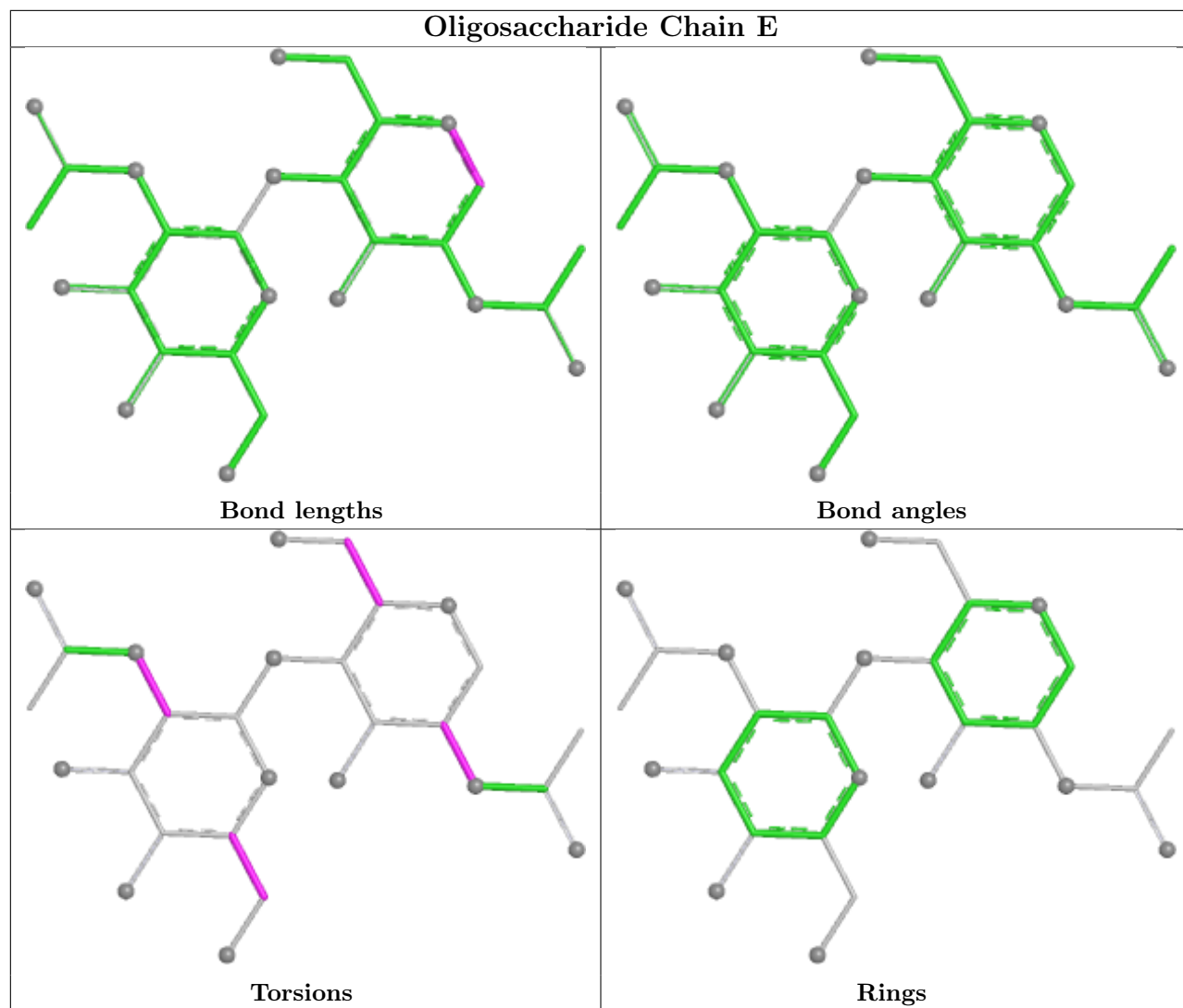
There are no ring outliers.

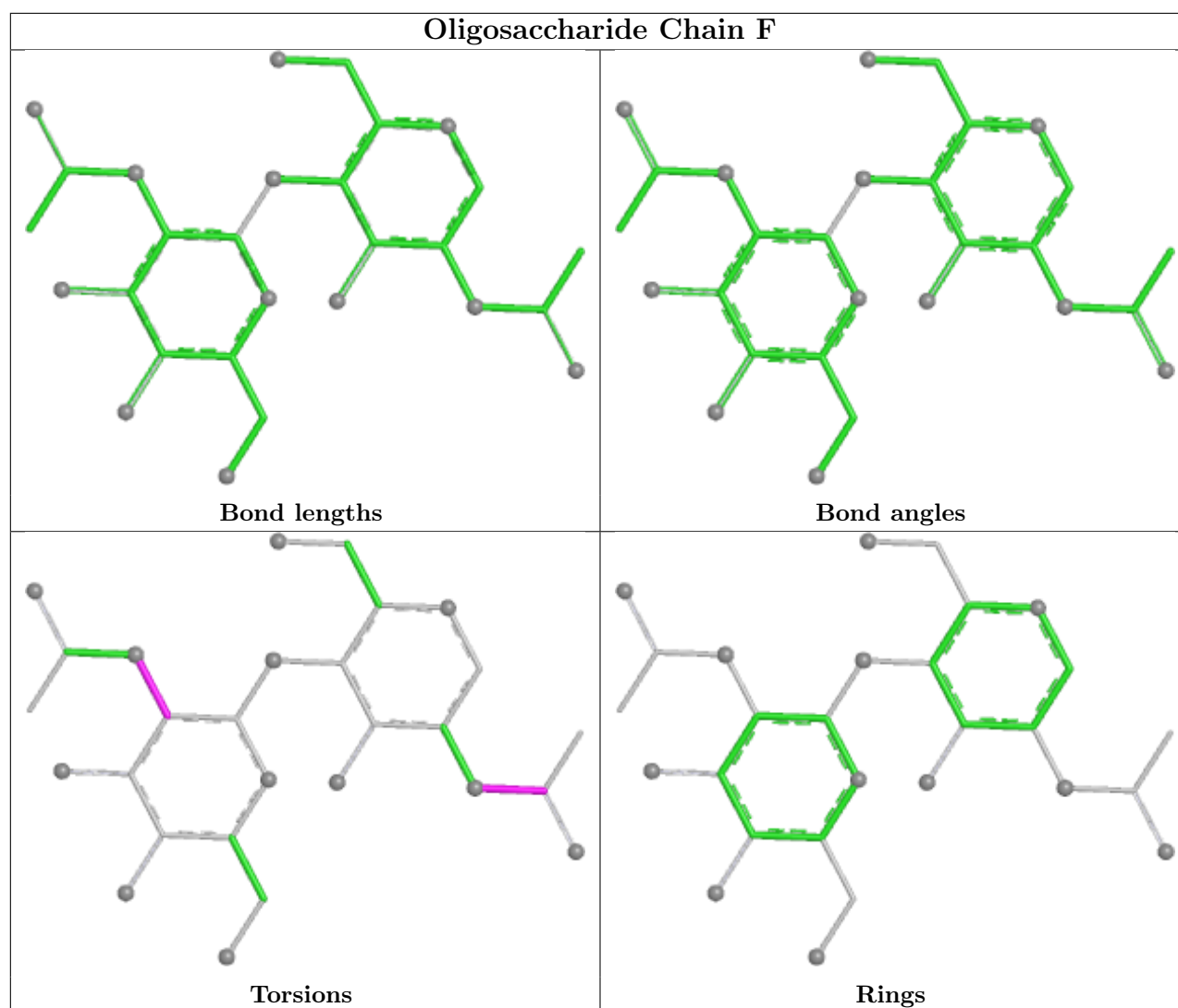
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	2	NAG	1	0
4	F	1	NAG	2	0
4	E	1	NAG	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](#)

18 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	1307	3	14,14,15	0.87	1 (7%)	17,19,21	2.53	3 (17%)
5	NAG	C	1303	3	14,14,15	0.23	0	17,19,21	0.37	0
5	NAG	C	1304	3	14,14,15	0.27	0	17,19,21	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	B	1304	3	14,14,15	0.34	0	17,19,21	0.39	0
5	NAG	A	1302	3	14,14,15	0.44	0	17,19,21	0.80	1 (5%)
5	NAG	A	1306	3	14,14,15	0.18	0	17,19,21	0.55	0
5	NAG	A	1309	3	14,14,15	0.22	0	17,19,21	0.49	0
5	NAG	B	1305	3	14,14,15	0.19	0	17,19,21	0.55	0
5	NAG	A	1301	3	14,14,15	0.23	0	17,19,21	0.49	0
5	NAG	B	1302	3	14,14,15	0.30	0	17,19,21	0.57	0
5	NAG	C	1302	3	14,14,15	0.24	0	17,19,21	0.48	0
5	NAG	B	1301	3	14,14,15	0.20	0	17,19,21	0.41	0
5	NAG	C	1301	3	14,14,15	0.20	0	17,19,21	0.50	0
5	NAG	B	1303	3	14,14,15	0.25	0	17,19,21	0.56	0
5	NAG	A	1305	3	14,14,15	0.65	1 (7%)	17,19,21	0.54	0
5	NAG	A	1304	3	14,14,15	0.22	0	17,19,21	0.46	0
5	NAG	A	1308	3	14,14,15	0.80	1 (7%)	17,19,21	1.19	2 (11%)
5	NAG	A	1303	3	14,14,15	0.16	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1307	3	-	3/6/23/26	0/1/1/1
5	NAG	C	1303	3	-	0/6/23/26	0/1/1/1
5	NAG	C	1304	3	-	2/6/23/26	0/1/1/1
5	NAG	B	1304	3	-	4/6/23/26	0/1/1/1
5	NAG	A	1302	3	-	4/6/23/26	0/1/1/1
5	NAG	A	1306	3	-	4/6/23/26	0/1/1/1
5	NAG	A	1309	3	-	3/6/23/26	0/1/1/1
5	NAG	B	1305	3	-	4/6/23/26	0/1/1/1
5	NAG	A	1301	3	-	4/6/23/26	0/1/1/1
5	NAG	B	1302	3	-	4/6/23/26	0/1/1/1
5	NAG	C	1302	3	-	1/6/23/26	0/1/1/1
5	NAG	B	1301	3	-	2/6/23/26	0/1/1/1
5	NAG	C	1301	3	-	2/6/23/26	0/1/1/1
5	NAG	B	1303	3	-	2/6/23/26	0/1/1/1
5	NAG	A	1305	3	-	1/6/23/26	0/1/1/1
5	NAG	A	1304	3	-	0/6/23/26	0/1/1/1
5	NAG	A	1308	3	-	3/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1303	3	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1307	NAG	O5-C1	-2.65	1.39	1.43
5	A	1305	NAG	O5-C1	-2.37	1.39	1.43
5	A	1308	NAG	O5-C1	2.03	1.47	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1307	NAG	C2-N2-C7	8.66	134.50	122.90
5	A	1307	NAG	C1-C2-N2	-4.10	103.97	110.43
5	A	1308	NAG	C1-C2-N2	-3.12	105.52	110.43
5	A	1308	NAG	C2-N2-C7	3.03	126.96	122.90
5	A	1302	NAG	C1-O5-C5	2.64	115.72	112.19
5	A	1307	NAG	C1-O5-C5	2.16	115.08	112.19

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1307	NAG	C1-C2-N2-C7
5	A	1307	NAG	C8-C7-N2-C2
5	A	1307	NAG	O7-C7-N2-C2
5	B	1303	NAG	C4-C5-C6-O6
5	B	1302	NAG	O5-C5-C6-O6
5	B	1304	NAG	C4-C5-C6-O6
5	A	1301	NAG	C4-C5-C6-O6
5	B	1303	NAG	O5-C5-C6-O6
5	A	1301	NAG	O5-C5-C6-O6
5	A	1302	NAG	O5-C5-C6-O6
5	C	1304	NAG	O5-C5-C6-O6
5	A	1302	NAG	C4-C5-C6-O6
5	C	1304	NAG	C4-C5-C6-O6
5	B	1302	NAG	C4-C5-C6-O6
5	B	1304	NAG	O5-C5-C6-O6
5	A	1301	NAG	C8-C7-N2-C2
5	A	1301	NAG	O7-C7-N2-C2
5	A	1302	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
5	A	1302	NAG	O7-C7-N2-C2
5	B	1304	NAG	C8-C7-N2-C2
5	B	1304	NAG	O7-C7-N2-C2
5	B	1301	NAG	O5-C5-C6-O6
5	B	1305	NAG	C4-C5-C6-O6
5	C	1302	NAG	O5-C5-C6-O6
5	B	1305	NAG	O5-C5-C6-O6
5	A	1308	NAG	O7-C7-N2-C2
5	A	1305	NAG	O5-C5-C6-O6
5	A	1303	NAG	C4-C5-C6-O6
5	A	1309	NAG	O5-C5-C6-O6
5	A	1308	NAG	C1-C2-N2-C7
5	A	1306	NAG	C4-C5-C6-O6
5	C	1301	NAG	C4-C5-C6-O6
5	C	1301	NAG	O5-C5-C6-O6
5	A	1306	NAG	C3-C2-N2-C7
5	A	1309	NAG	C3-C2-N2-C7
5	B	1302	NAG	C3-C2-N2-C7
5	B	1305	NAG	C3-C2-N2-C7
5	A	1303	NAG	O5-C5-C6-O6
5	A	1306	NAG	C1-C2-N2-C7
5	A	1309	NAG	C1-C2-N2-C7
5	B	1302	NAG	C1-C2-N2-C7
5	B	1305	NAG	C1-C2-N2-C7
5	A	1308	NAG	C3-C2-N2-C7
5	A	1306	NAG	O5-C5-C6-O6
5	B	1301	NAG	C4-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1307	NAG	3	0
5	B	1304	NAG	1	0
5	A	1309	NAG	1	0
5	A	1308	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

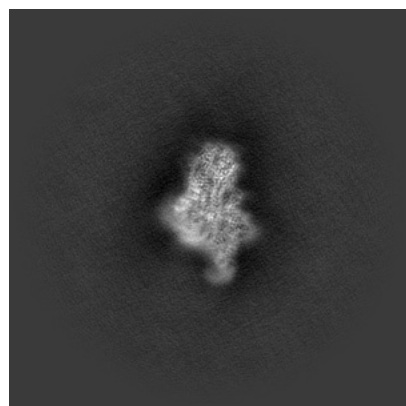
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-28756. These allow visual inspection of the internal detail of the map and identification of artifacts.

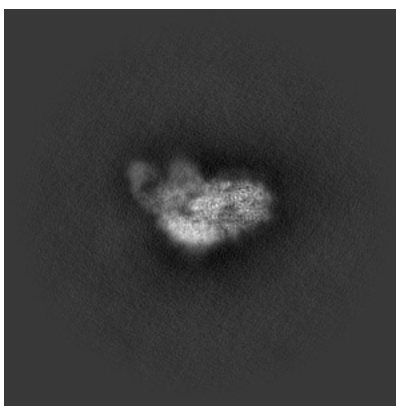
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

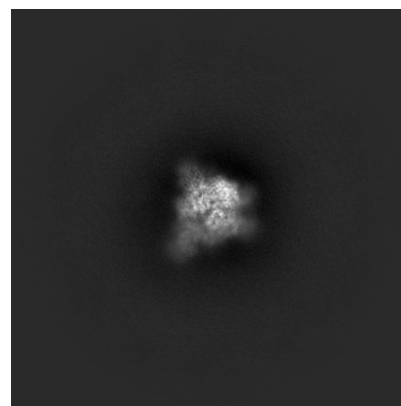
6.1.1 Primary map



X

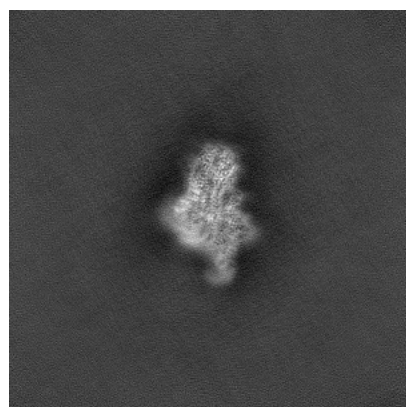


Y

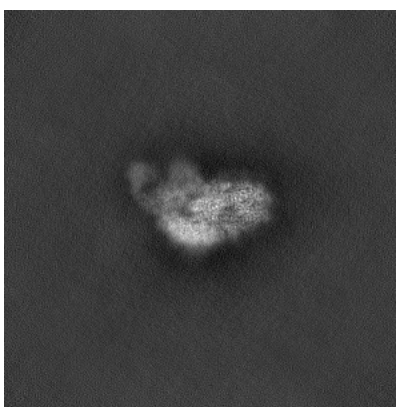


Z

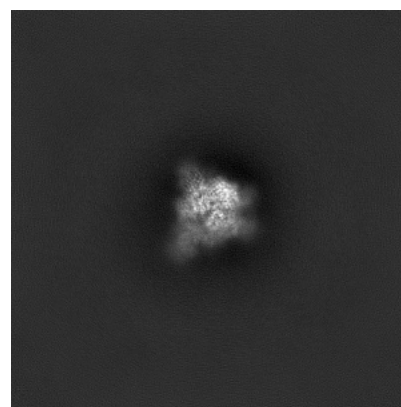
6.1.2 Raw map



X



Y

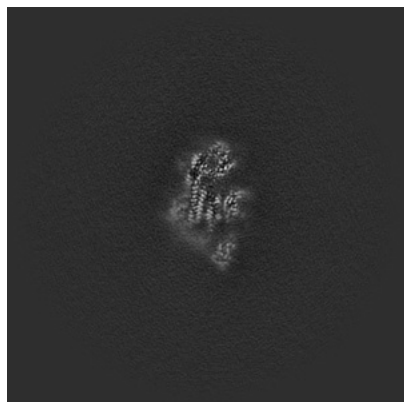


Z

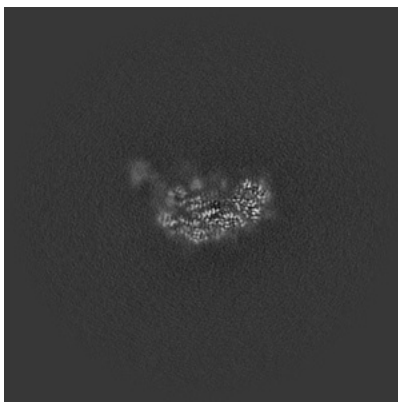
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

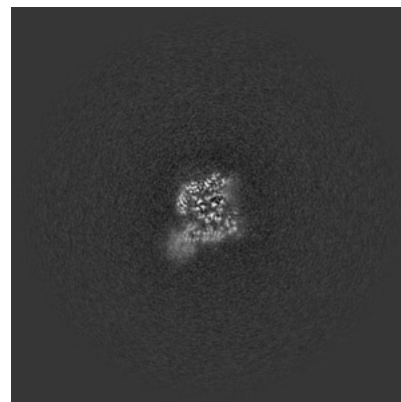
6.2.1 Primary map



X Index: 200

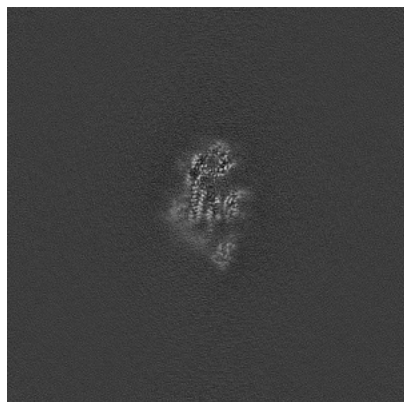


Y Index: 200

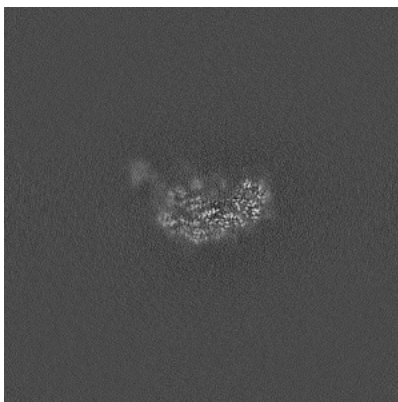


Z Index: 200

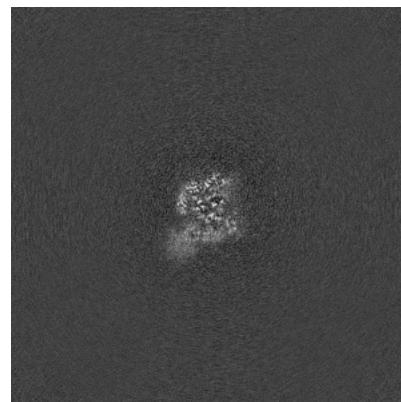
6.2.2 Raw map



X Index: 200



Y Index: 200

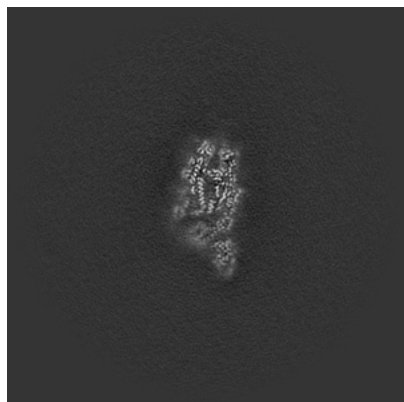


Z Index: 200

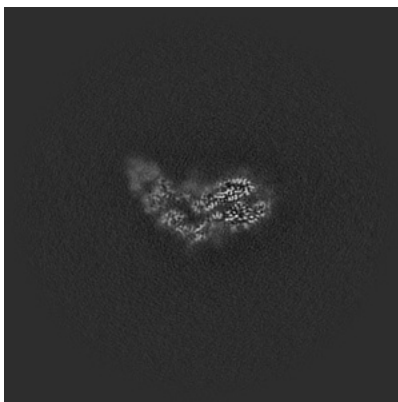
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

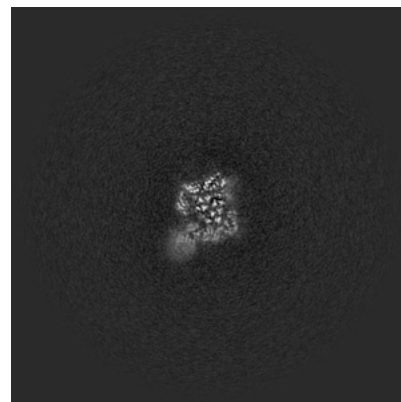
6.3.1 Primary map



X Index: 210

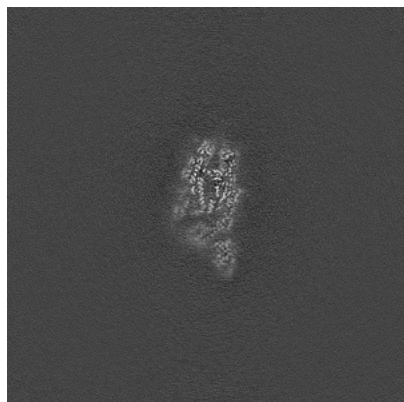


Y Index: 210

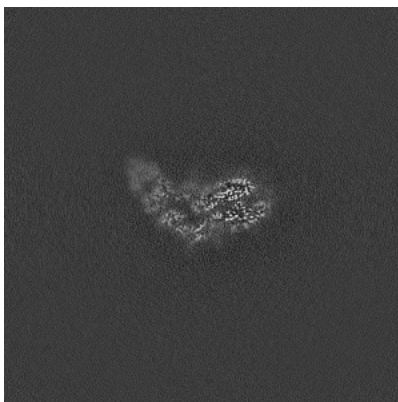


Z Index: 201

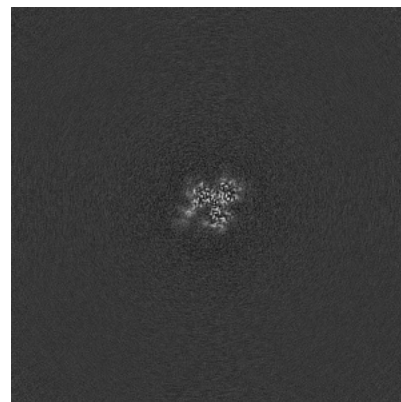
6.3.2 Raw map



X Index: 210



Y Index: 210

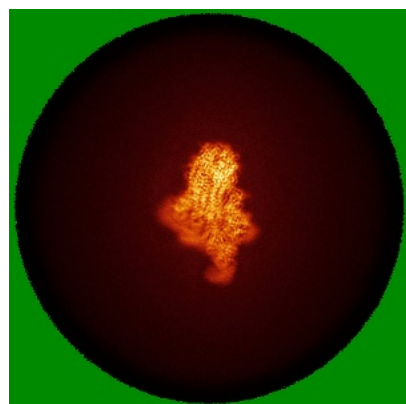


Z Index: 231

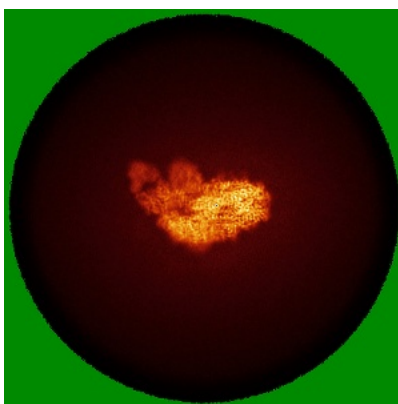
The images above show the largest variance slices of the map in three orthogonal directions.

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

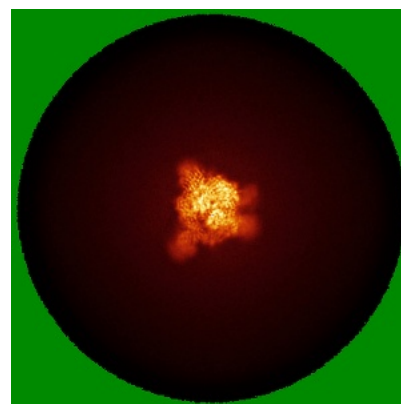
6.4.1 Primary map



X

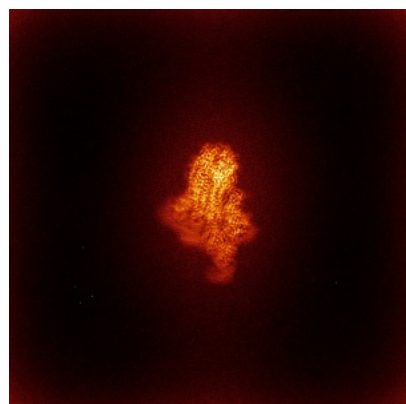


Y

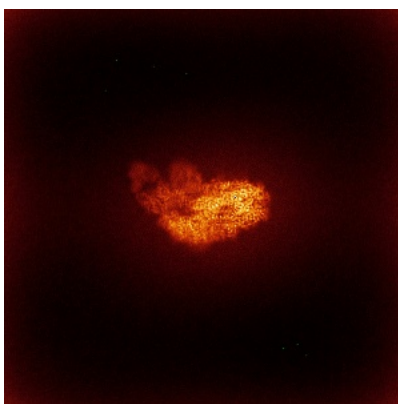


Z

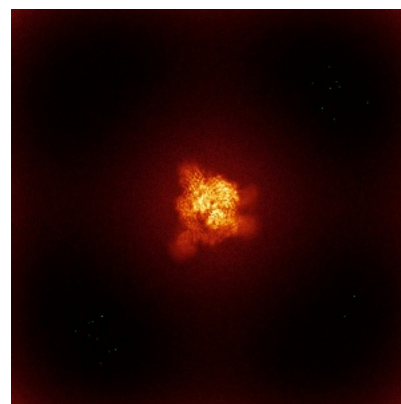
6.4.2 Raw map



X



Y

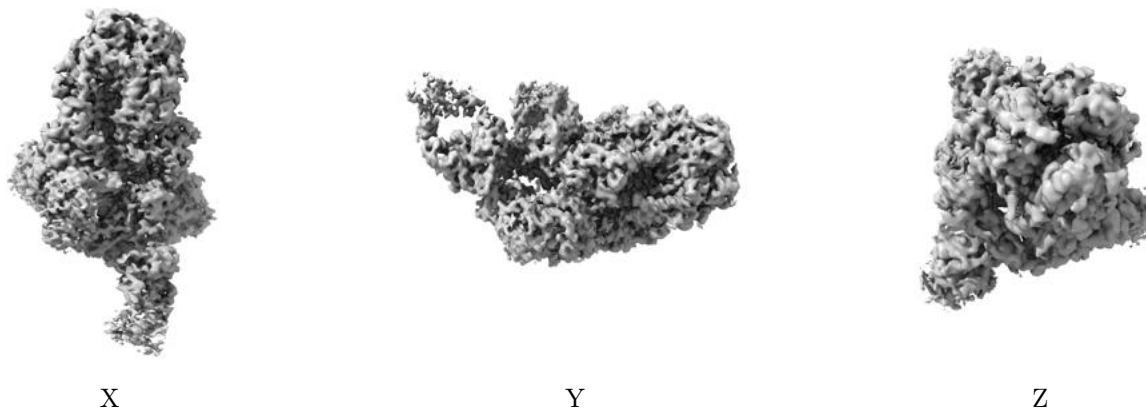


Z

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.

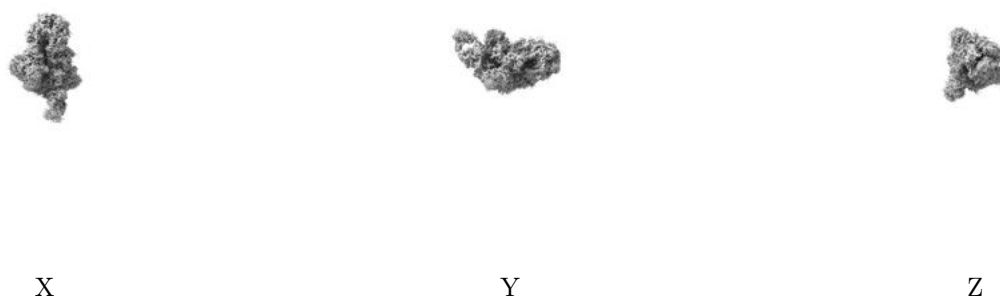
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.2. 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.

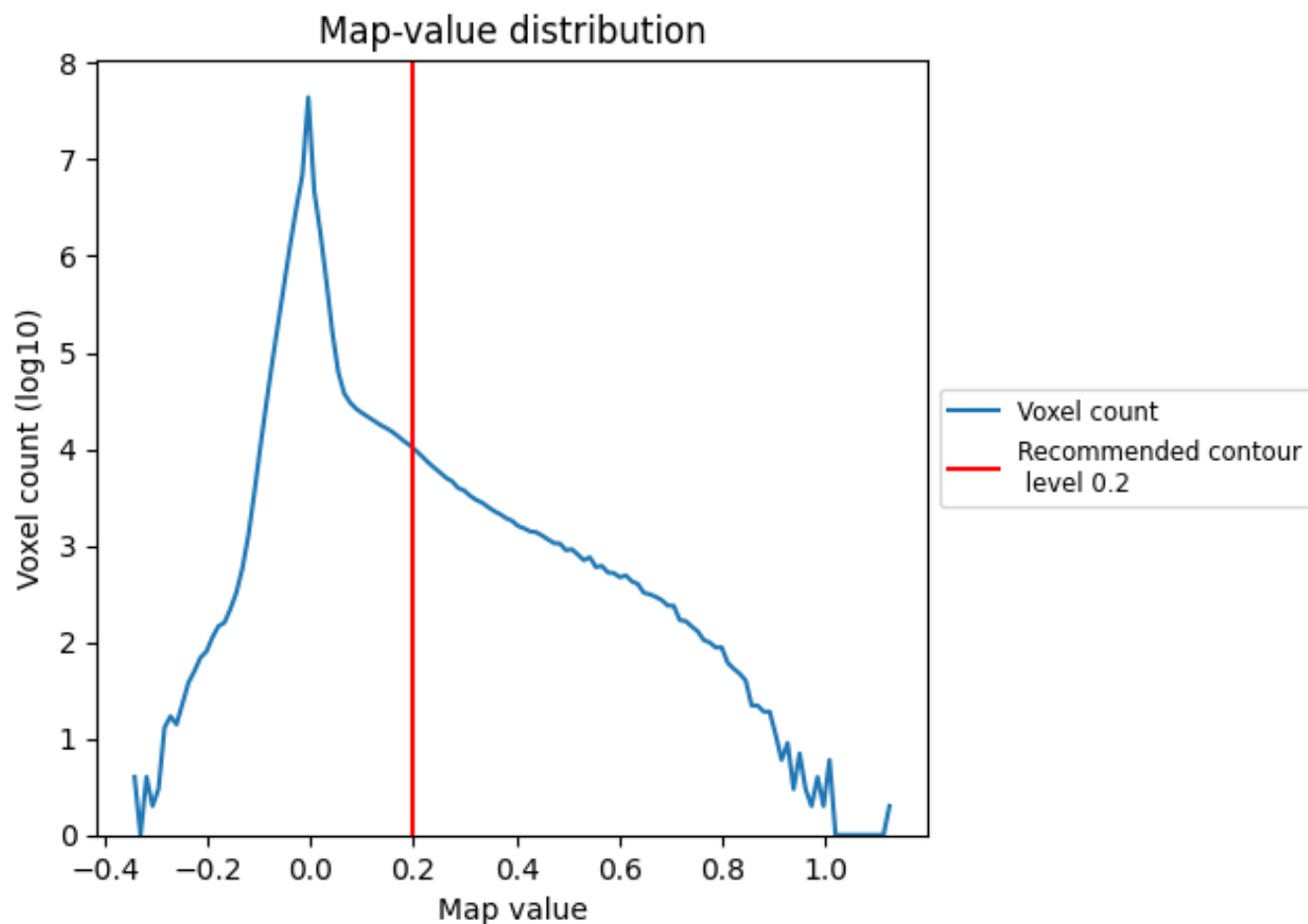
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

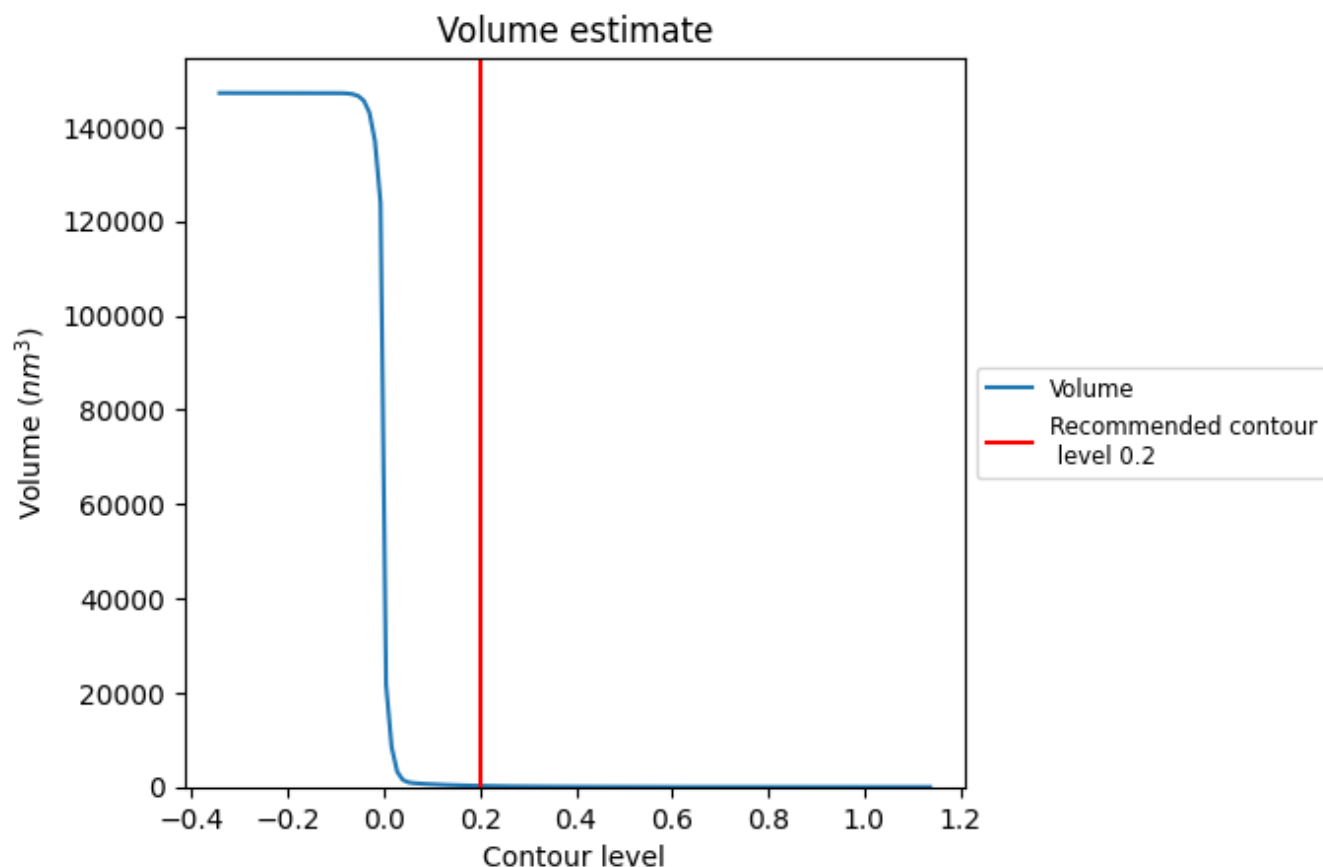
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

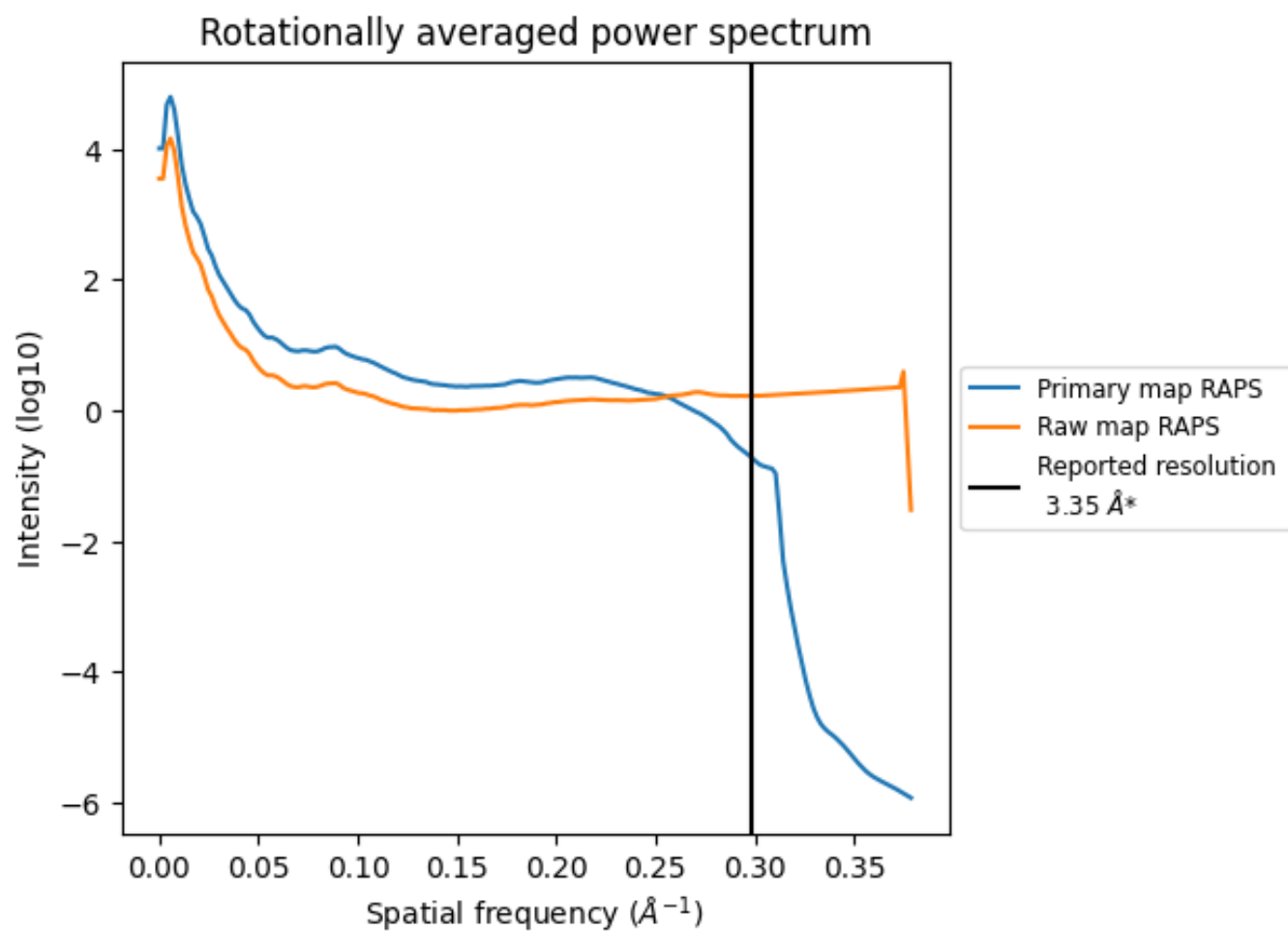
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 232 nm^3 ; this corresponds to an approximate mass of 210 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 [i](#)

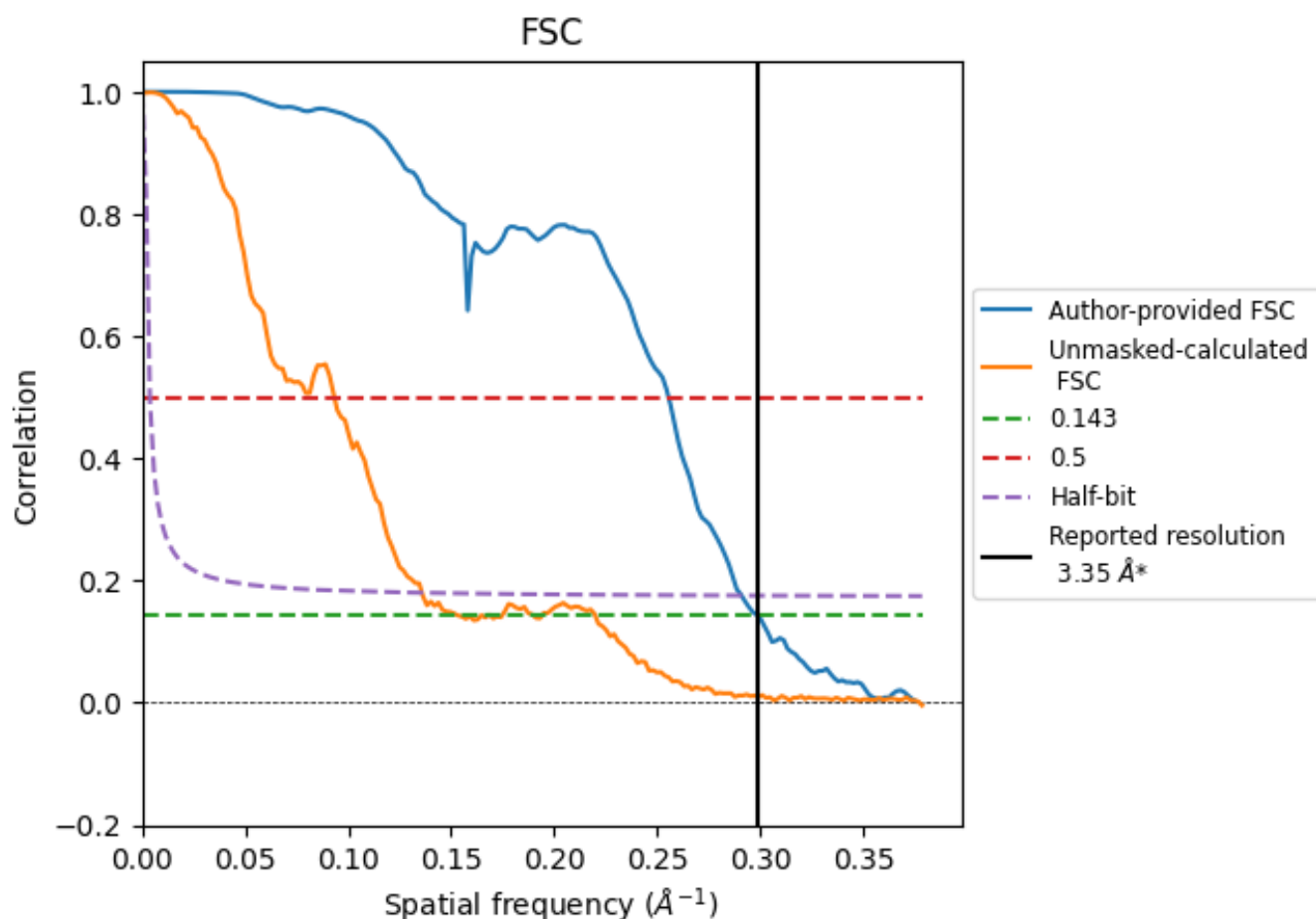


*Reported resolution corresponds to spatial frequency of $0.299 \text{ \AA}^{-1}$

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 0.299 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

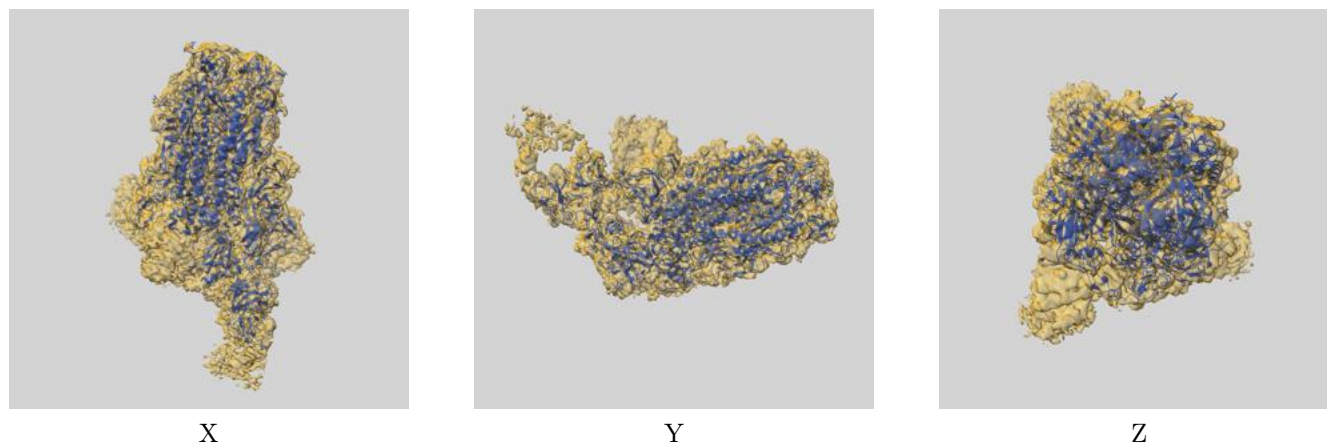
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.35	-	-
Author-provided FSC curve	3.35	3.91	3.44
Unmasked-calculated*	6.53	10.72	7.36

*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 6.53 differs from the reported value 3.35 by more than 10 %

9 Map-model fit [i](#)

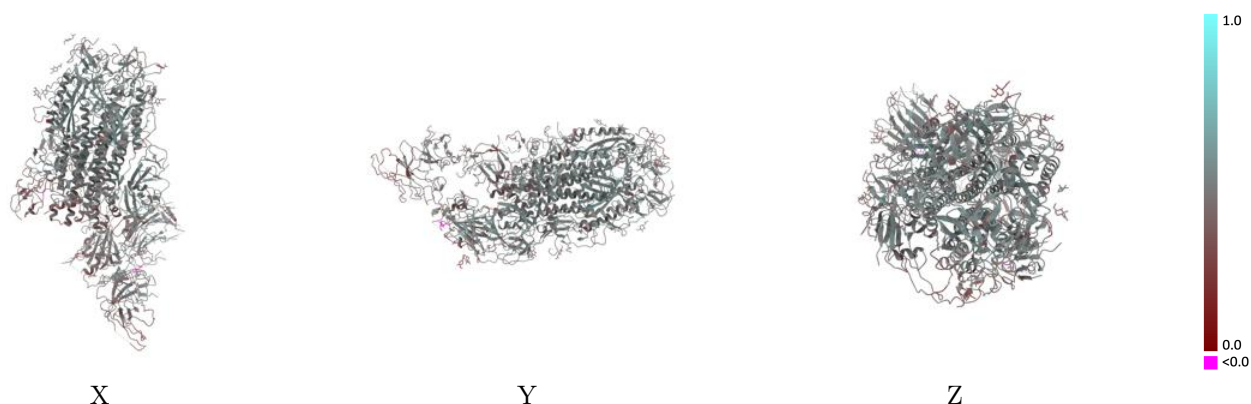
This section contains information regarding the fit between EMDB map EMD-28756 and PDB model 8F0G. Per-residue inclusion information can be found in section [3](#) on page [12](#).

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.2 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](#)

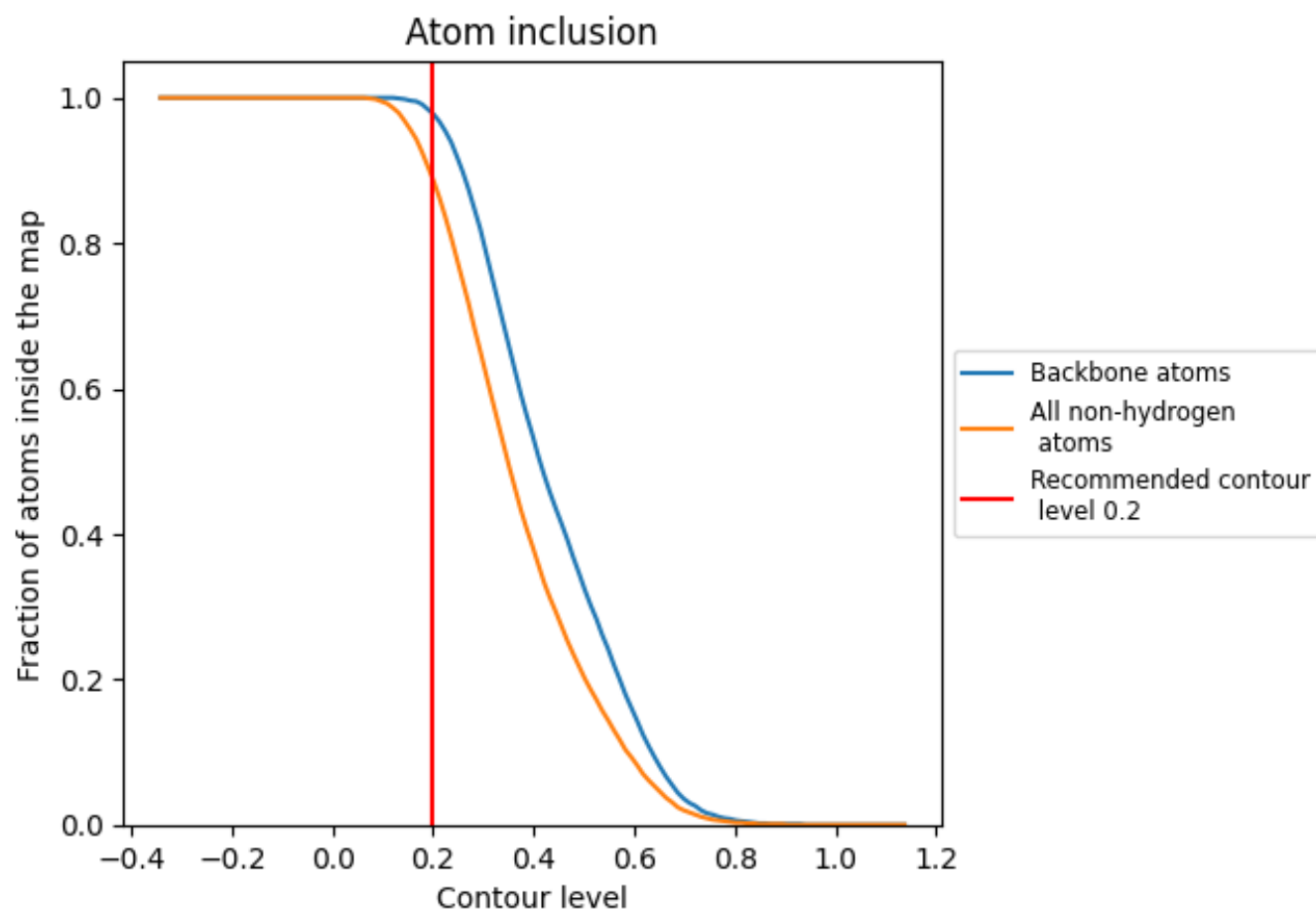


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 89% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8870	0.4690
A	0.9040	0.4840
B	0.8890	0.4630
C	0.8990	0.4730
D	0.6430	0.2930
E	0.7500	0.3900
F	0.6790	0.4130
X	0.7500	0.3920
Y	0.8000	0.4500

1.0

0.0

<0.0