



Full wwPDB EM Validation Report ⓘ

Apr 11, 2023 – 10:31 am BST

PDB ID : 8AQW
EMDB ID : EMD-15592
Title : BA.4/5 SARS-CoV-2 Spike bound to mouse ACE2 (local)
Authors : Lau, K.; Ni, D.; Beckert, B.; Nazarov, S.; Myasnikov, A.; Pojer, F.; Stahlberg, H.; Uchikawa, E.
Deposited on : 2022-08-13
Resolution : 3.30 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.dev50
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.32.2

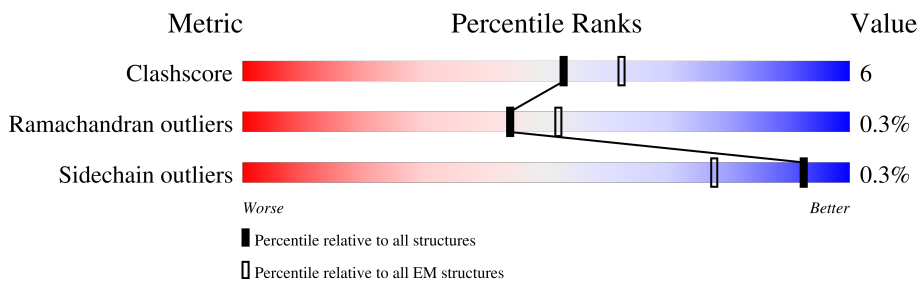
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.30 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	B	1283	
2	A	884	
3	C	2	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6457 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Spike glycoprotein, Fibritin.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	B	193	1543	994	261	280	8	0	0

There are 95 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	24	ILE	THR	variant	UNP P0DTC2
B	?	-	LEU	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	?	-	PRO	deletion	UNP P0DTC2
B	29	SER	ALA	variant	UNP P0DTC2
B	?	-	HIS	deletion	UNP P0DTC2
B	?	-	VAL	deletion	UNP P0DTC2
B	142	ASP	GLY	variant	UNP P0DTC2
B	213	GLY	VAL	variant	UNP P0DTC2
B	339	ASP	GLY	variant	UNP P0DTC2
B	371	PHE	SER	variant	UNP P0DTC2
B	373	PRO	SER	variant	UNP P0DTC2
B	375	PHE	SER	variant	UNP P0DTC2
B	376	ALA	THR	variant	UNP P0DTC2
B	405	ASN	ASP	variant	UNP P0DTC2
B	408	SER	ARG	variant	UNP P0DTC2
B	417	ASN	LYS	variant	UNP P0DTC2
B	440	LYS	ASN	variant	UNP P0DTC2
B	452	ARG	LEU	variant	UNP P0DTC2
B	477	ASN	SER	variant	UNP P0DTC2
B	478	LYS	THR	variant	UNP P0DTC2
B	484	ALA	GLU	variant	UNP P0DTC2
B	486	VAL	PHE	variant	UNP P0DTC2
B	498	ARG	GLN	variant	UNP P0DTC2
B	501	TYR	ASN	variant	UNP P0DTC2
B	505	HIS	TYR	variant	UNP P0DTC2
B	614	GLY	ASP	variant	UNP P0DTC2
B	655	TYR	HIS	variant	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	679	LYS	ASN	variant	UNP P0DTC2
B	681	HIS	PRO	variant	UNP P0DTC2
B	682	GLY	ARG	variant	UNP P0DTC2
B	683	SER	ARG	variant	UNP P0DTC2
B	685	SER	ARG	variant	UNP P0DTC2
B	764	LYS	ASN	variant	UNP P0DTC2
B	796	TYR	ASP	variant	UNP P0DTC2
B	954	HIS	GLN	variant	UNP P0DTC2
B	969	LYS	ASN	variant	UNP P0DTC2
B	986	PRO	LYS	variant	UNP P0DTC2
B	987	PRO	VAL	variant	UNP P0DTC2
B	1209	GLY	-	linker	UNP P0DTC2
B	1210	SER	-	linker	UNP P0DTC2
B	1211	GLY	-	linker	UNP P0DTC2
B	1212	TYR	-	linker	UNP P0DTC2
B	1232	LEU	PHE	engineered mutation	UNP P10104
B	1238	GLY	-	expression tag	UNP P10104
B	1239	ARG	-	expression tag	UNP P10104
B	1240	SER	-	expression tag	UNP P10104
B	1241	LEU	-	expression tag	UNP P10104
B	1242	GLU	-	expression tag	UNP P10104
B	1243	VAL	-	expression tag	UNP P10104
B	1244	LEU	-	expression tag	UNP P10104
B	1245	PHE	-	expression tag	UNP P10104
B	1246	GLN	-	expression tag	UNP P10104
B	1247	GLY	-	expression tag	UNP P10104
B	1248	PRO	-	expression tag	UNP P10104
B	1249	GLY	-	expression tag	UNP P10104
B	1250	HIS	-	expression tag	UNP P10104
B	1251	HIS	-	expression tag	UNP P10104
B	1252	HIS	-	expression tag	UNP P10104
B	1253	HIS	-	expression tag	UNP P10104
B	1254	HIS	-	expression tag	UNP P10104
B	1255	HIS	-	expression tag	UNP P10104
B	1256	HIS	-	expression tag	UNP P10104
B	1257	HIS	-	expression tag	UNP P10104
B	1258	SER	-	expression tag	UNP P10104
B	1259	ALA	-	expression tag	UNP P10104
B	1260	TRP	-	expression tag	UNP P10104
B	1261	SER	-	expression tag	UNP P10104
B	1262	HIS	-	expression tag	UNP P10104
B	1263	PRO	-	expression tag	UNP P10104

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1264	GLN	-	expression tag	UNP P10104
B	1265	PHE	-	expression tag	UNP P10104
B	1266	GLU	-	expression tag	UNP P10104
B	1267	LYS	-	expression tag	UNP P10104
B	1268	GLY	-	expression tag	UNP P10104
B	1269	GLY	-	expression tag	UNP P10104
B	1270	GLY	-	expression tag	UNP P10104
B	1271	SER	-	expression tag	UNP P10104
B	1272	GLY	-	expression tag	UNP P10104
B	1273	GLY	-	expression tag	UNP P10104
B	1274	GLY	-	expression tag	UNP P10104
B	1275	GLY	-	expression tag	UNP P10104
B	1276	SER	-	expression tag	UNP P10104
B	1277	GLY	-	expression tag	UNP P10104
B	1278	GLY	-	expression tag	UNP P10104
B	1279	SER	-	expression tag	UNP P10104
B	1280	ALA	-	expression tag	UNP P10104
B	1281	TRP	-	expression tag	UNP P10104
B	1282	SER	-	expression tag	UNP P10104
B	1283	HIS	-	expression tag	UNP P10104
B	1284	PRO	-	expression tag	UNP P10104
B	1285	GLN	-	expression tag	UNP P10104
B	1286	PHE	-	expression tag	UNP P10104
B	1287	GLU	-	expression tag	UNP P10104
B	1288	LYS	-	expression tag	UNP P10104

- Molecule 2 is a protein called Processed angiotensin-converting enzyme 2,Ig gamma-2A chain C region, A allele.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	A	596	4872	3103	819	921	29	1	0

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	initiating methionine	UNP Q8R0I0
A	-14	GLY	-	expression tag	UNP Q8R0I0
A	-13	THR	-	expression tag	UNP Q8R0I0
A	-12	LEU	-	expression tag	UNP Q8R0I0
A	-11	SER	-	expression tag	UNP Q8R0I0
A	-10	ALA	-	expression tag	UNP Q8R0I0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	PRO	-	expression tag	UNP Q8R0I0
A	-8	PRO	-	expression tag	UNP Q8R0I0
A	-7	CYS	-	expression tag	UNP Q8R0I0
A	-6	THR	-	expression tag	UNP Q8R0I0
A	-5	GLN	-	expression tag	UNP Q8R0I0
A	-4	ARG	-	expression tag	UNP Q8R0I0
A	-3	ILE	-	expression tag	UNP Q8R0I0
A	-2	LYS	-	expression tag	UNP Q8R0I0
A	-1	TRP	-	expression tag	UNP Q8R0I0
A	0	LYS	-	expression tag	UNP Q8R0I0
A	1	GLY	-	expression tag	UNP Q8R0I0
A	2	LEU	-	expression tag	UNP Q8R0I0
A	3	LEU	-	expression tag	UNP Q8R0I0
A	4	LEU	-	expression tag	UNP Q8R0I0
A	5	THR	-	expression tag	UNP Q8R0I0
A	6	ALA	-	expression tag	UNP Q8R0I0
A	7	SER	-	expression tag	UNP Q8R0I0
A	8	LEU	-	expression tag	UNP Q8R0I0
A	9	LEU	-	expression tag	UNP Q8R0I0
A	10	ASN	-	expression tag	UNP Q8R0I0
A	11	PHE	-	expression tag	UNP Q8R0I0
A	12	TRP	-	expression tag	UNP Q8R0I0
A	13	ASN	-	expression tag	UNP Q8R0I0
A	14	LEU	-	expression tag	UNP Q8R0I0
A	15	PRO	-	expression tag	UNP Q8R0I0
A	16	THR	-	expression tag	UNP Q8R0I0
A	17	THR	-	expression tag	UNP Q8R0I0
A	18	ALA	-	expression tag	UNP Q8R0I0
A	616	LEU	-	linker	UNP Q8R0I0
A	617	GLU	-	linker	UNP Q8R0I0
A	618	VAL	-	linker	UNP Q8R0I0
A	619	LEU	-	linker	UNP Q8R0I0
A	620	PHE	-	linker	UNP Q8R0I0
A	621	GLN	-	linker	UNP Q8R0I0
A	622	GLY	-	linker	UNP Q8R0I0
A	623	PRO	-	linker	UNP Q8R0I0
A	624	MET	-	linker	UNP Q8R0I0
A	625	ASP	-	linker	UNP Q8R0I0
A	859	HIS	-	expression tag	UNP P01863
A	860	HIS	-	expression tag	UNP P01863
A	861	HIS	-	expression tag	UNP P01863
A	862	HIS	-	expression tag	UNP P01863

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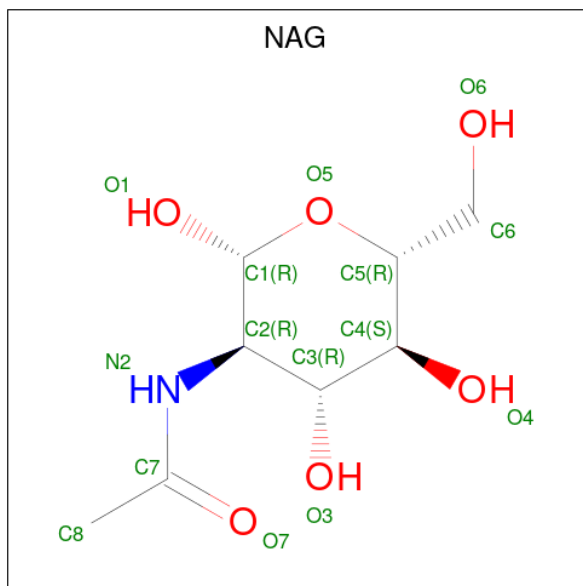
Chain	Residue	Modelled	Actual	Comment	Reference
A	863	HIS	-	expression tag	UNP P01863
A	864	HIS	-	expression tag	UNP P01863
A	865	HIS	-	expression tag	UNP P01863
A	866	HIS	-	expression tag	UNP P01863
A	867	HIS	-	expression tag	UNP P01863
A	868	HIS	-	expression tag	UNP P01863

- Molecule 3 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	
			Total	C	N			O
3	C	2	28	16	2	10	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).

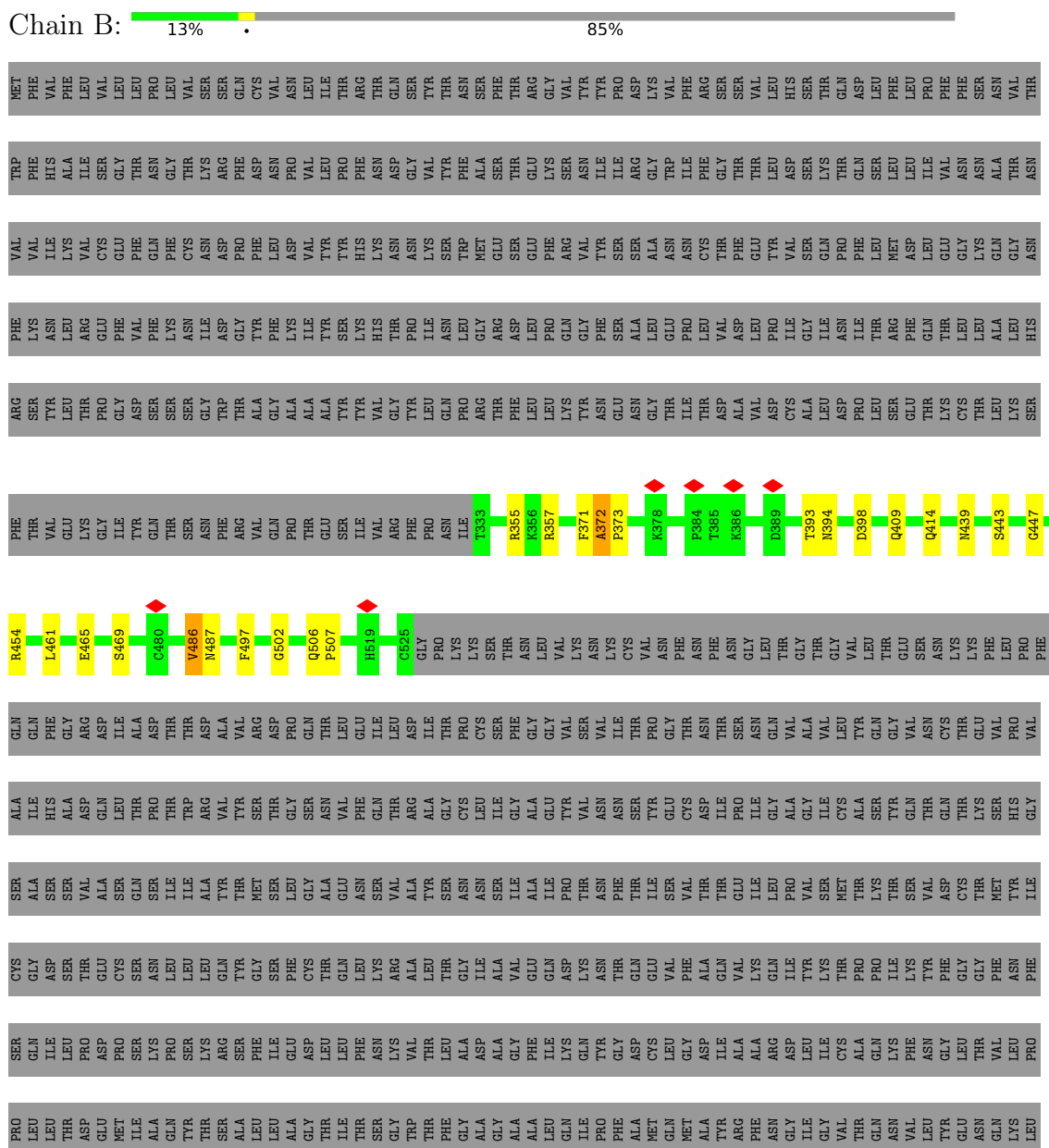


Mol	Chain	Residues	Atoms			AltConf	
			Total	C	N		O
4	B	1	14	8	1	5	0

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: Spike glycoprotein,Fibrinin



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

• Molecule 2: Processed angiotensin-converting enzyme 2,Ig gamma-2A chain C region, A allele



MET	GLY	THR	THR	SER	ALA	PRO	PRO	CYS	THR	THR	GLN	ARG	ILE	LYS	TRP	GLY	LEU	LEU	LEU	THR	THR
Q89	P90	P91	P92	R95	Q96	L97	Q98	S103	G104	S105	A110	K114	Q115	L116	M117	T118	L119	L120	M121	T122	T125
L162	W163	G167	W168	R169	G173	K174	R177	P178	Y183	M190	R191	R192	A193	N194	R204	A212	D213	N218	R219	N220	A231
T276	N277	T282	V283	P284	A286	Q287	K288	P289	D292	V293	T294	M297	F315	L320	E335	D338	C361	T365	H493	D494	E495
E435	T445	E457	K458	M462	V463	F464	R465	G466	I468	K475	K476	M480	K481	R482	E483	I484	V488	P492	H493	D494	E495
F592	D593	W594	L595	K596	Q598	N599	R600	N601	S602	F603	V604	N607	T608	E609	W610	S611	A614	D615	LEU	GLU	VAL
PHE	PRO	ALA	PRO	LYS	ILE	LYS	ASP	VAL	GLN	ASP	MET	VAL	ILE	VAL	THR	THR	CYS	VAL	VAL	VAL	VAL
VAL	SER	ALA	LEU	PRO	ILE	LYS	ASP	GLN	HIS	VAL	THR	ASP	GLN	VAL	VAL	VAL	VAL	VAL	VAL	VAL	VAL
VAL	THR	LEU	THR	CYS	MET	VAL	THR	ASP	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR

TYR
SER
CYS
SER
VAL
VAL
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GLU
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- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:

100%

WAG1
WAG2

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	103496	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	58	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2800	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	2.114	Depositor
Minimum map value	-0.745	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.172	Depositor
Map size (\AA)	365.184, 365.184, 365.184	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.0144, 1.0144, 1.0144	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	B	0.26	0/1589	0.52	0/2163
2	A	0.24	0/5010	0.45	0/6801
All	All	0.25	0/6599	0.46	0/8964

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	B	1543	0	1466	16	0
2	A	4872	0	4639	55	0
3	C	28	0	25	0	0
4	B	14	0	13	0	0
All	All	6457	0	6143	71	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (71) 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)
2:A:85:LEU:HD11	2:A:97:LEU:HD23	1.69	0.72
2:A:34[C]:GLN:OE1	2:A:35:GLU:N	2.23	0.70
2:A:458:LYS:O	2:A:462:MET:HG2	1.95	0.66
1:B:393:THR:HG22	1:B:394:ASN:H	1.61	0.66
2:A:419:LYS:NZ	2:A:428:PHE:O	2.29	0.65
2:A:247:LYS:HG2	2:A:282:THR:HA	1.79	0.65
1:B:439:ASN:O	1:B:443:SER:OG	2.18	0.62
1:B:461:LEU:HD22	1:B:465:GLU:HB3	1.81	0.61
1:B:454:ARG:NH2	1:B:469:SER:O	2.35	0.60
2:A:245:ARG:NH1	2:A:260:GLY:O	2.33	0.60
2:A:553:LYS:HA	2:A:556:LYS:HE3	1.84	0.59
2:A:190:MET:SD	2:A:194:ASN:ND2	2.75	0.59
2:A:288:LYS:HE2	2:A:289:PRO:HD2	1.84	0.59
2:A:294:THR:HG23	2:A:365:THR:HA	1.84	0.59
2:A:177:ARG:NH1	2:A:495:GLU:O	2.31	0.58
2:A:249:MET:HA	2:A:256:ILE:HG21	1.87	0.57
2:A:494:ASP:OD1	2:A:495:GLU:N	2.38	0.56
2:A:177:ARG:NH1	2:A:497:TYR:O	2.40	0.55
2:A:135:PRO:HD3	2:A:163:TRP:NE1	2.22	0.55
2:A:457:GLU:HG2	2:A:513:ILE:HD13	1.89	0.55
1:B:355:ARG:HA	1:B:398:ASP:HB3	1.89	0.54
2:A:92:ILE:HD12	2:A:92:ILE:H	1.73	0.54
2:A:276:THR:HG1	2:A:445:THR:HG1	1.56	0.53
2:A:168:TRP:NE1	2:A:502:SER:OG	2.42	0.53
1:B:372:ALA:HB3	1:B:373:PRO:HD3	1.92	0.52
1:B:439:ASN:HB2	1:B:506:GLN:HE21	1.75	0.52
2:A:607:ASN:ND2	2:A:609:GLU:OE1	2.40	0.52
2:A:37:GLU:OE2	2:A:393:ARG:NH1	2.40	0.52
2:A:592:PHE:HE1	2:A:596:LYS:HD2	1.74	0.51
2:A:263:PRO:HG2	2:A:266:LEU:HD12	1.93	0.50
1:B:439:ASN:CB	1:B:506:GLN:HE21	2.24	0.50
2:A:315:PHE:HD1	2:A:320:LEU:HD12	1.77	0.50
2:A:28:PHE:HE2	2:A:80:ALA:HB2	1.78	0.49
2:A:414:THR:HG23	2:A:417:HIS:H	1.78	0.48
1:B:486:VAL:HG12	1:B:487:ASN:N	2.29	0.48
1:B:502:GLY:O	1:B:506:GLN:HB2	2.14	0.47
2:A:378:HIS:CE1	2:A:401:HIS:HB3	2.49	0.47
2:A:406:GLU:HG3	2:A:518:ARG:NH2	2.29	0.47
2:A:91:PRO:O	2:A:95:ARG:HG3	2.15	0.46
2:A:387:ARG:HH22	2:A:559:SER:HB3	1.81	0.46
2:A:402:GLU:HB2	2:A:518:ARG:HB3	1.98	0.45
1:B:371:PHE:O	1:B:373:PRO:HD3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:204:ARG:HH22	2:A:465:ARG:HH12	1.66	0.44
2:A:374:HIS:ND1	2:A:406:GLU:OE2	2.46	0.44
2:A:88:ILE:HG22	2:A:89:GLN:N	2.32	0.44
2:A:480:MET:O	2:A:484:ILE:HB	2.18	0.44
2:A:78:LYS:HE3	2:A:78:LYS:HB2	1.74	0.44
2:A:417:HIS:HB2	2:A:543:ASP:OD2	2.19	0.43
2:A:544:ILE:HG13	2:A:544:ILE:O	2.18	0.43
2:A:396:ALA:HB3	2:A:400:PHE:CD2	2.54	0.43
2:A:83:PHE:O	2:A:85:LEU:HD12	2.19	0.43
2:A:232:GLU:OE1	2:A:232:GLU:N	2.48	0.43
2:A:592:PHE:CE1	2:A:596:LYS:HD2	2.54	0.43
2:A:406:GLU:HG3	2:A:518:ARG:HH21	1.84	0.42
2:A:55:THR:OG1	2:A:58:ASN:OD1	2.36	0.42
1:B:497:PHE:CG	1:B:507:PRO:HG3	2.54	0.42
2:A:120:LEU:HD23	2:A:120:LEU:HA	1.87	0.42
1:B:357:ARG:HE	1:B:394:ASN:HD22	1.67	0.42
2:A:284:PRO:HB3	2:A:594:TRP:HH2	1.85	0.41
2:A:218:ASN:OD1	2:A:219:ARG:N	2.53	0.41
2:A:285:PHE:CE2	2:A:433:GLU:HG3	2.56	0.41
2:A:482:ARG:HE	2:A:488:VAL:HG23	1.86	0.41
2:A:499:ASP:N	2:A:500:PRO:HD2	2.36	0.41
2:A:87:GLU:O	2:A:88:ILE:HG13	2.21	0.40
2:A:335:GLU:HG3	2:A:361:CYS:HB3	2.03	0.40
2:A:95:ARG:O	2:A:98:GLN:HG3	2.20	0.40
1:B:447:GLY:HA2	1:B:497:PHE:O	2.21	0.40
2:A:177:ARG:HB3	2:A:178:PRO:HD3	2.03	0.40
1:B:357:ARG:NE	1:B:394:ASN:HD22	2.20	0.40
2:A:68:LYS:HE2	2:A:68:LYS:HB3	1.79	0.40
1:B:409:GLN:HA	1:B:414:GLN:HG2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	191/1283 (15%)	174 (91%)	15 (8%)	2 (1%)	15	46
2	A	596/884 (67%)	560 (94%)	36 (6%)	0	100	100
All	All	787/2167 (36%)	734 (93%)	51 (6%)	2 (0%)	44	71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	372	ALA
1	B	486	VAL

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	166/1107 (15%)	166 (100%)	0	100	100
2	A	522/788 (66%)	518 (99%)	4 (1%)	81	89
All	All	688/1895 (36%)	684 (99%)	4 (1%)	92	91

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	34[A]	GLN
2	A	34[B]	GLN
2	A	34[C]	GLN
2	A	131	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	24	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

2 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
3	NAG	C	1	3,2	14,14,15	0.20	0	17,19,21	0.51	0
3	NAG	C	2	3	14,14,15	0.33	0	17,19,21	0.37	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	C	2	3	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2	NAG	C4-C5-C6-O6
3	C	1	NAG	O5-C5-C6-O6

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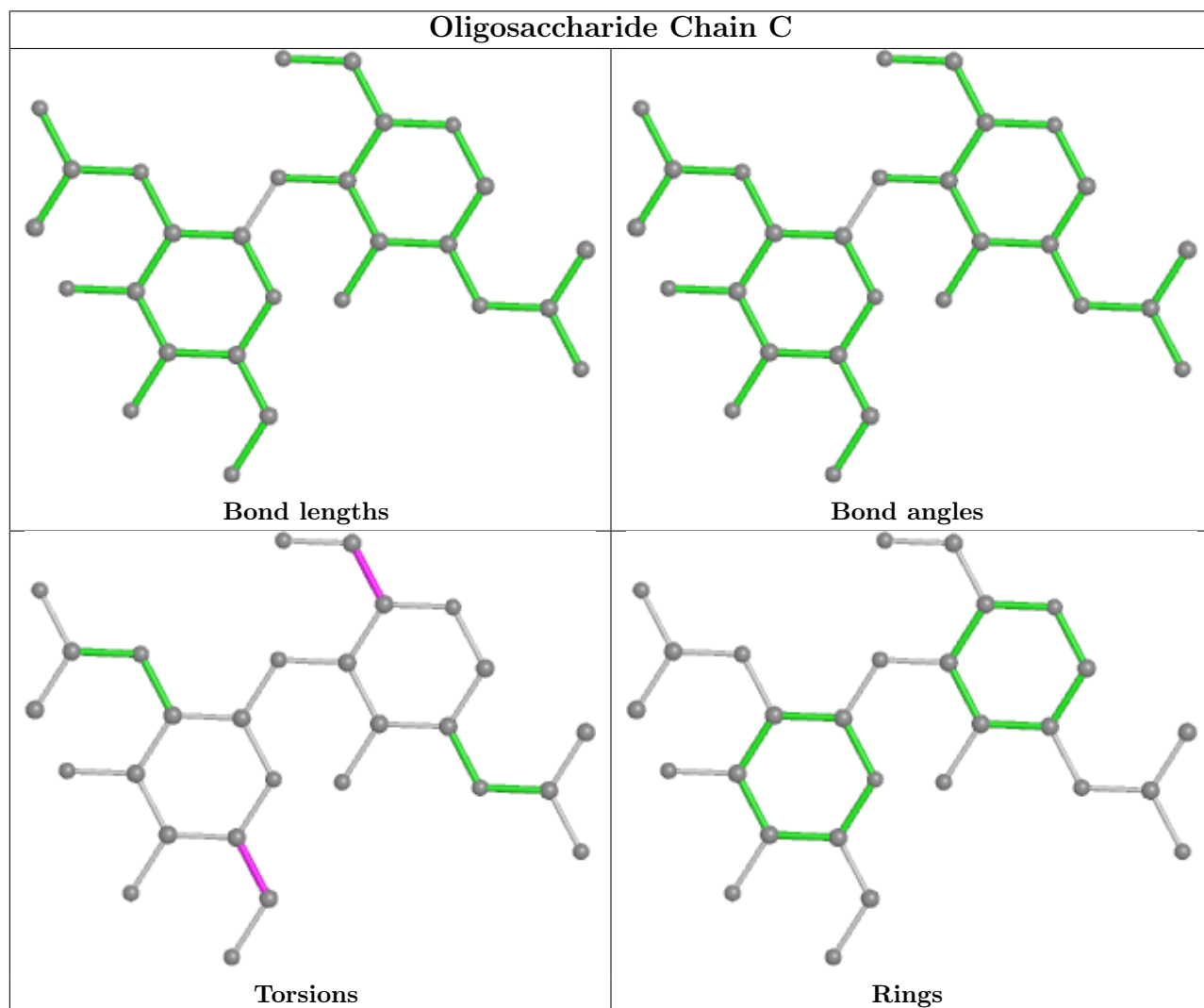
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Mol	Chain	Res	Type	Atoms
3	C	2	NAG	O5-C5-C6-O6
3	C	1	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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



5.6 Ligand geometry [i](#)

1 ligand is 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	B	1301	1	14,14,15	0.23	0	17,19,21	0.45	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	1301	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1301	NAG	O5-C5-C6-O6
4	B	1301	NAG	C4-C5-C6-O6
4	B	1301	NAG	C8-C7-N2-C2
4	B	1301	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

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.

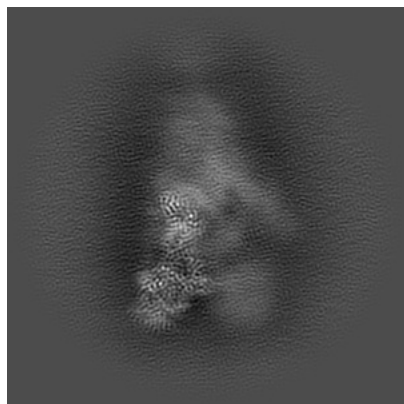
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-15592. These allow visual inspection of the internal detail of the map and identification of artifacts.

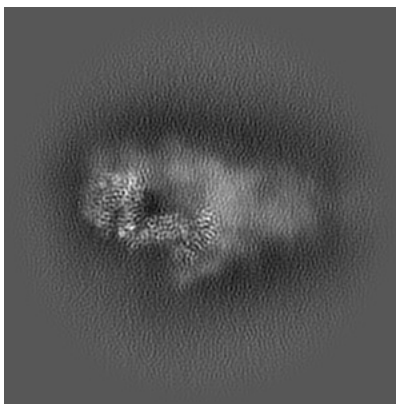
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

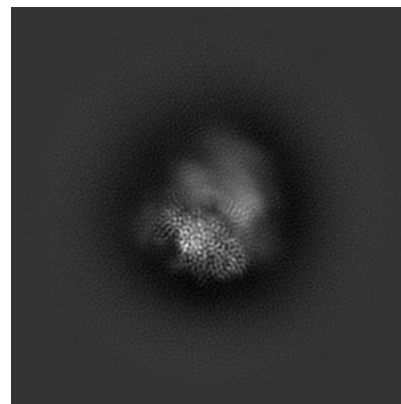
6.1.1 Primary map



X

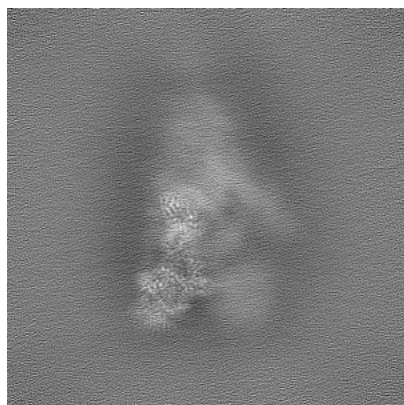


Y

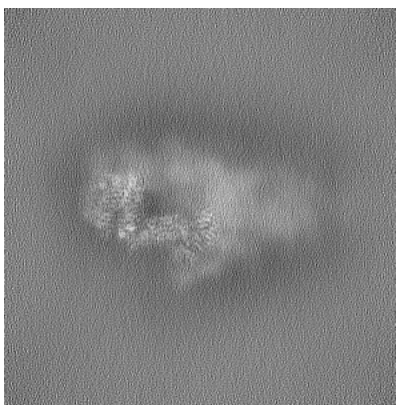


Z

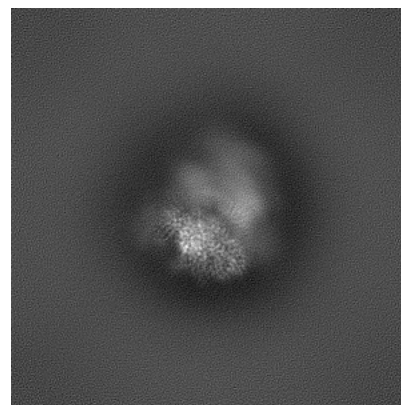
6.1.2 Raw map



X



Y

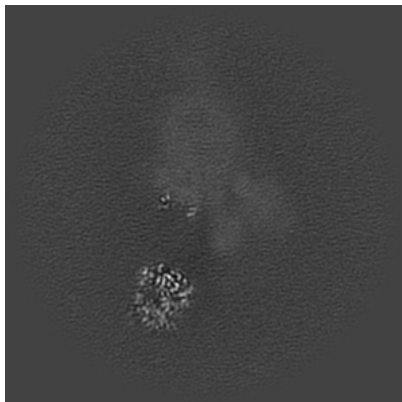


Z

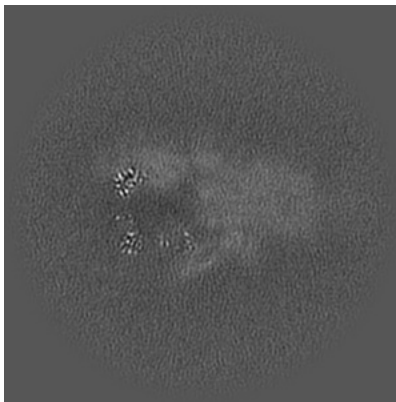
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

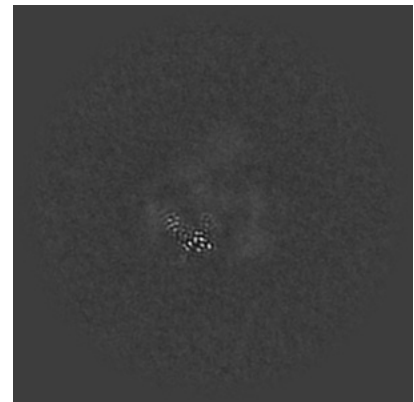
6.2.1 Primary map



X Index: 180

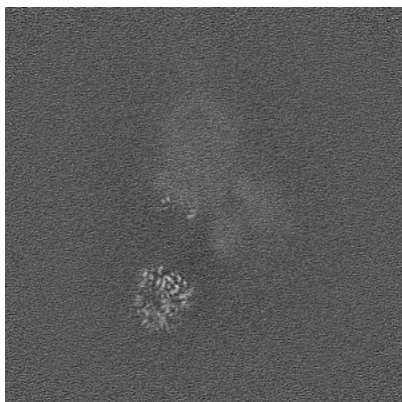


Y Index: 180

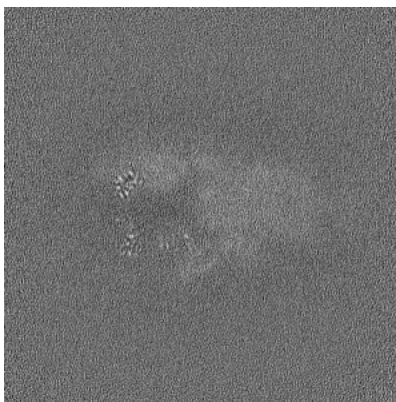


Z Index: 180

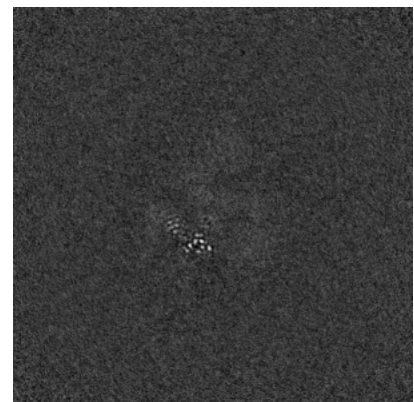
6.2.2 Raw map



X Index: 180



Y Index: 180

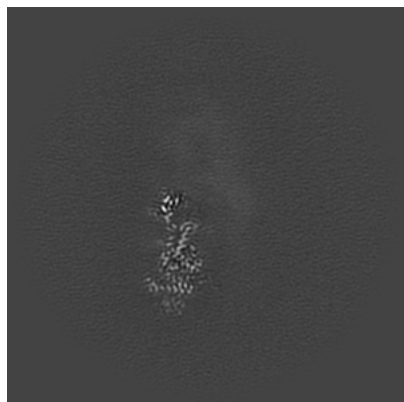


Z Index: 180

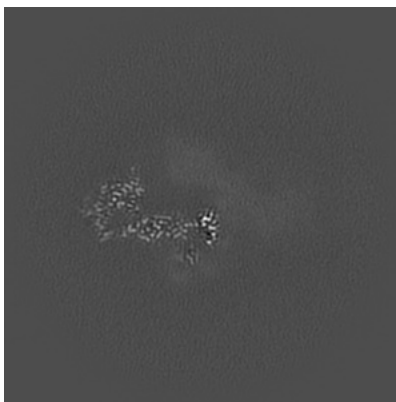
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

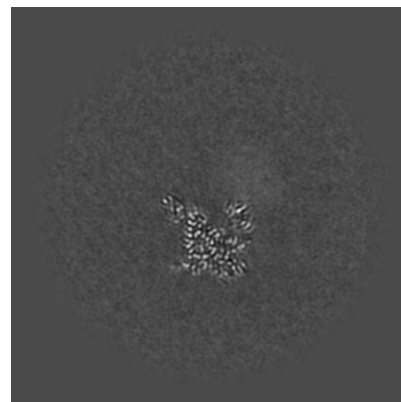
6.3.1 Primary map



X Index: 157

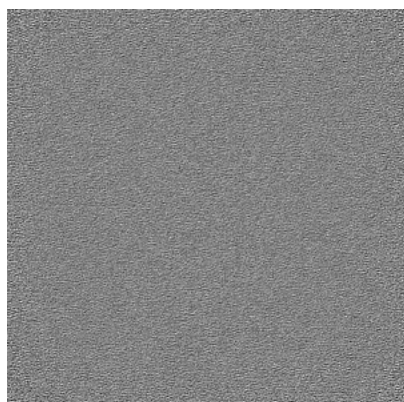


Y Index: 149

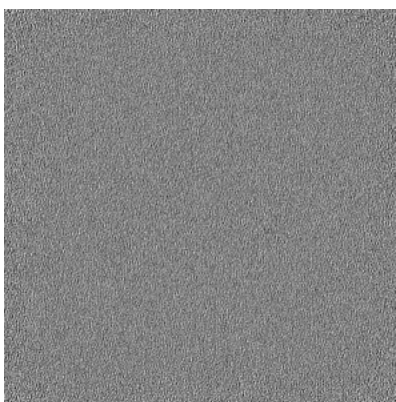


Z Index: 114

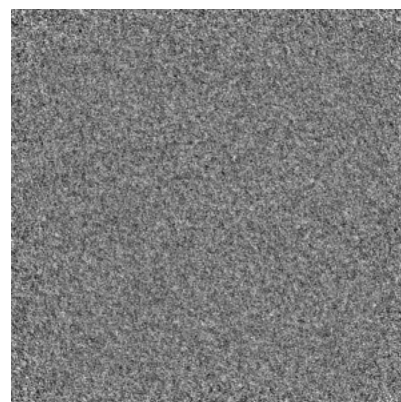
6.3.2 Raw map



X Index: 0



Y Index: 0

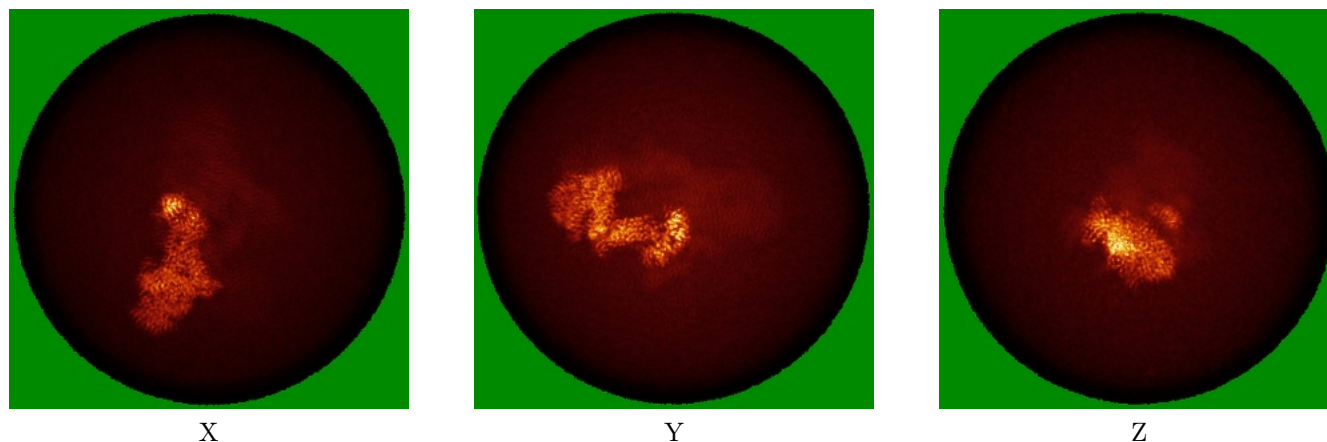


Z Index: 359

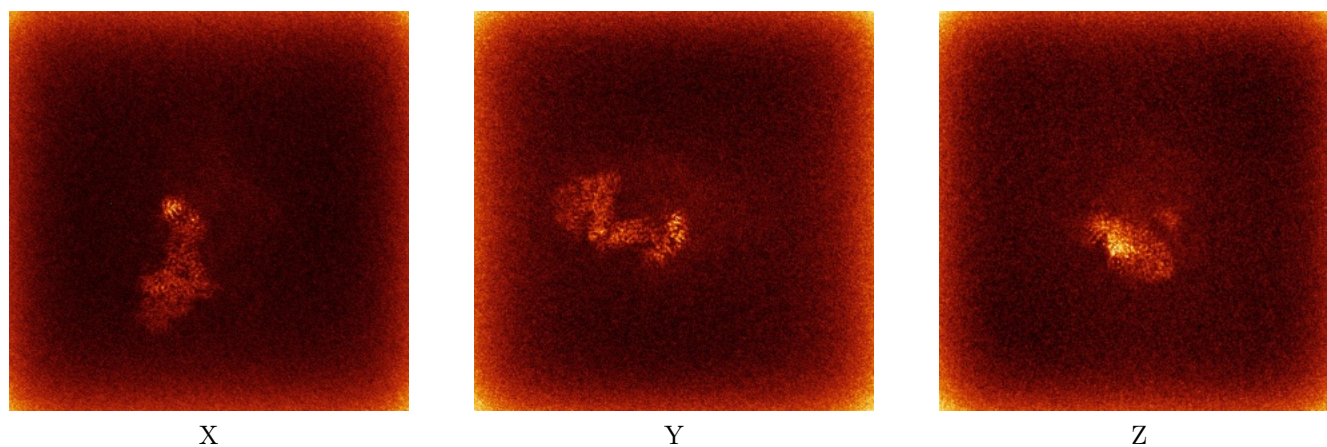
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



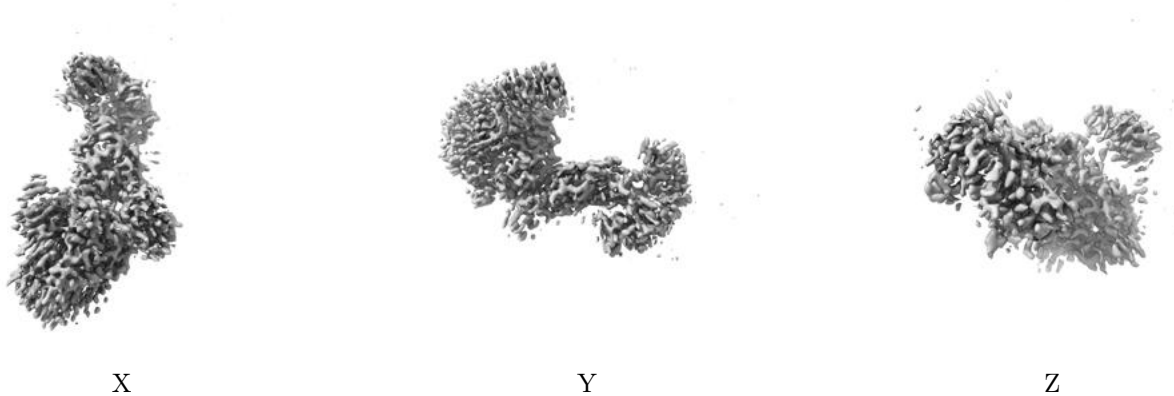
6.4.2 Raw map



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.

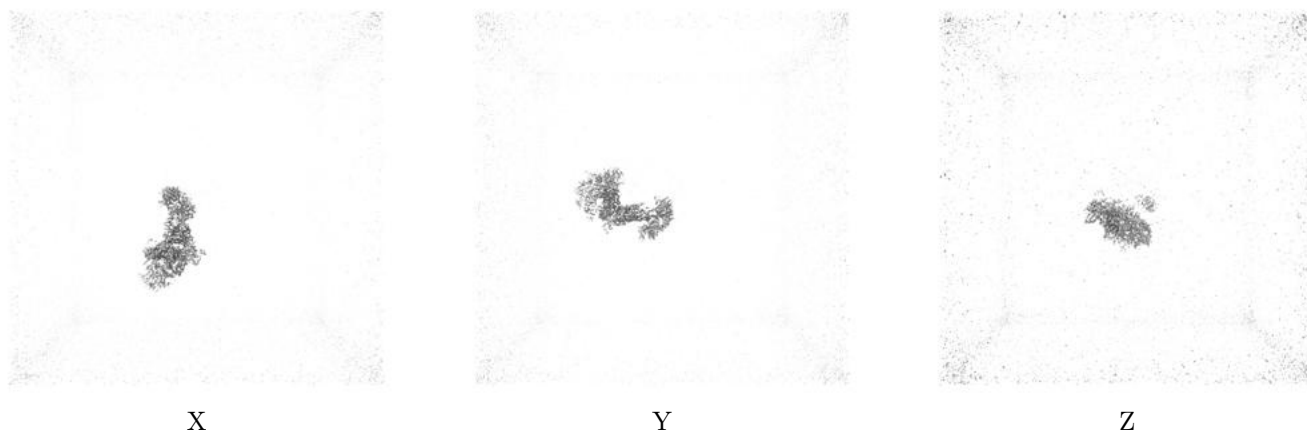
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.172. 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.

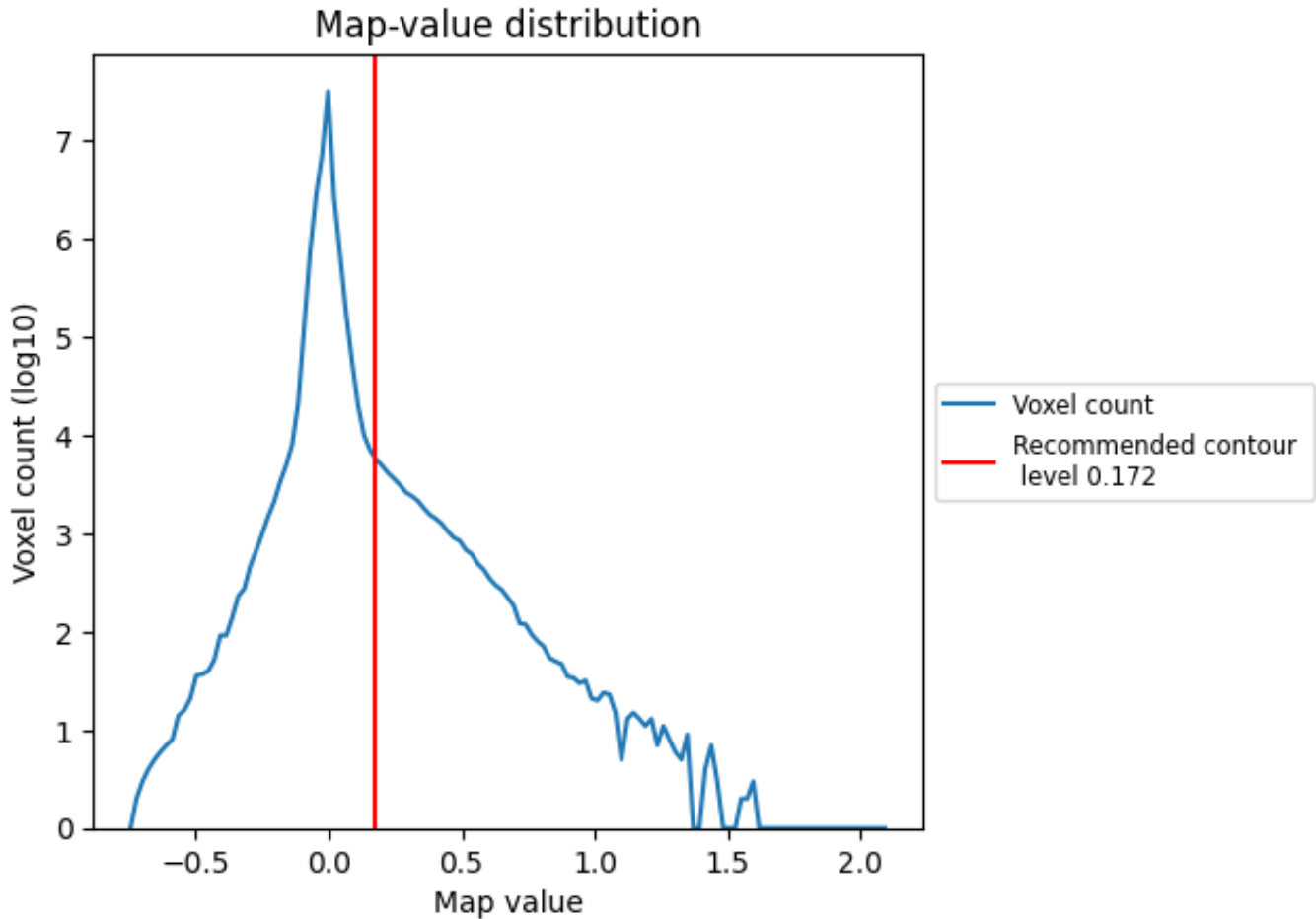
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

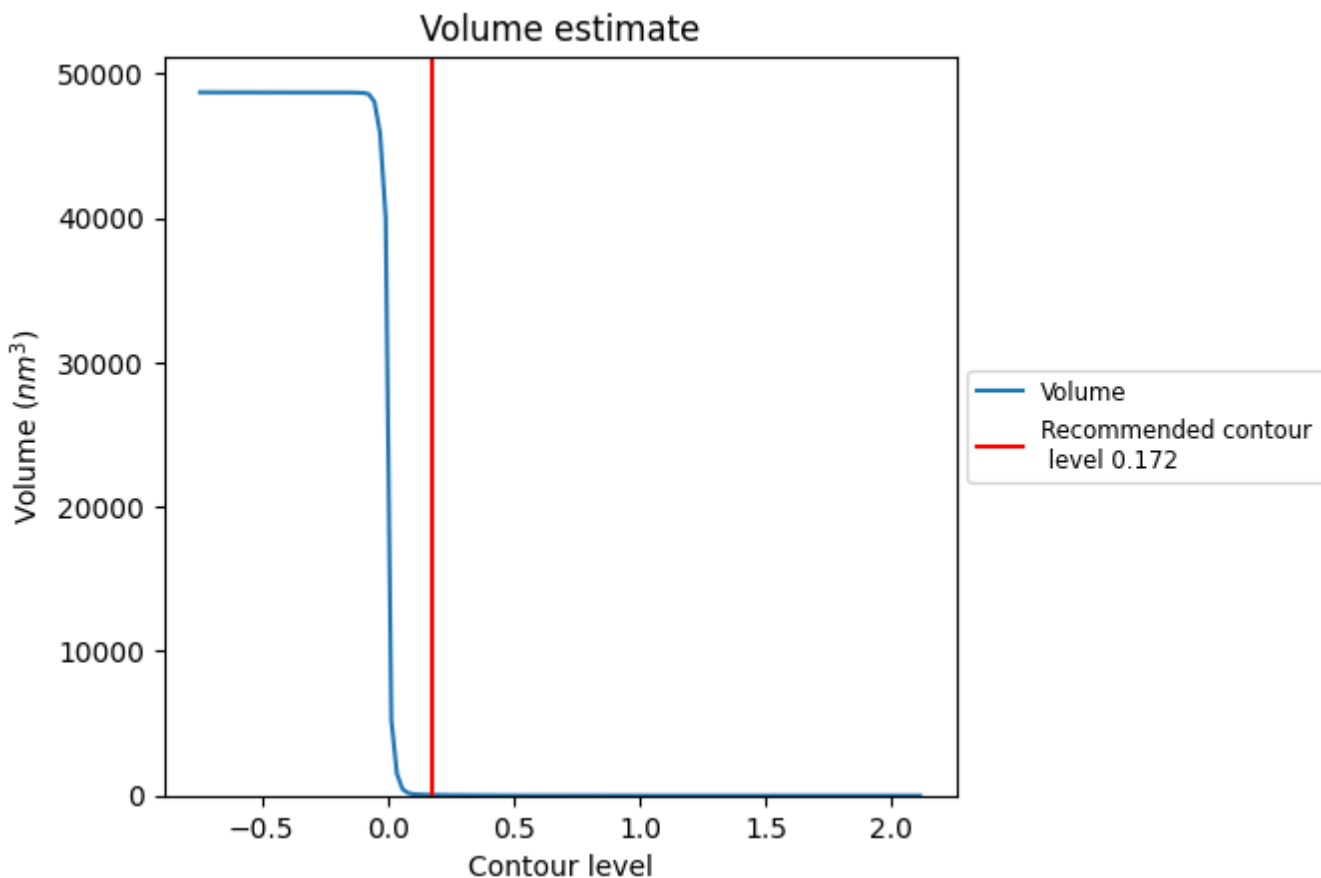
This section contains the results of statistical analysis of the map.

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.

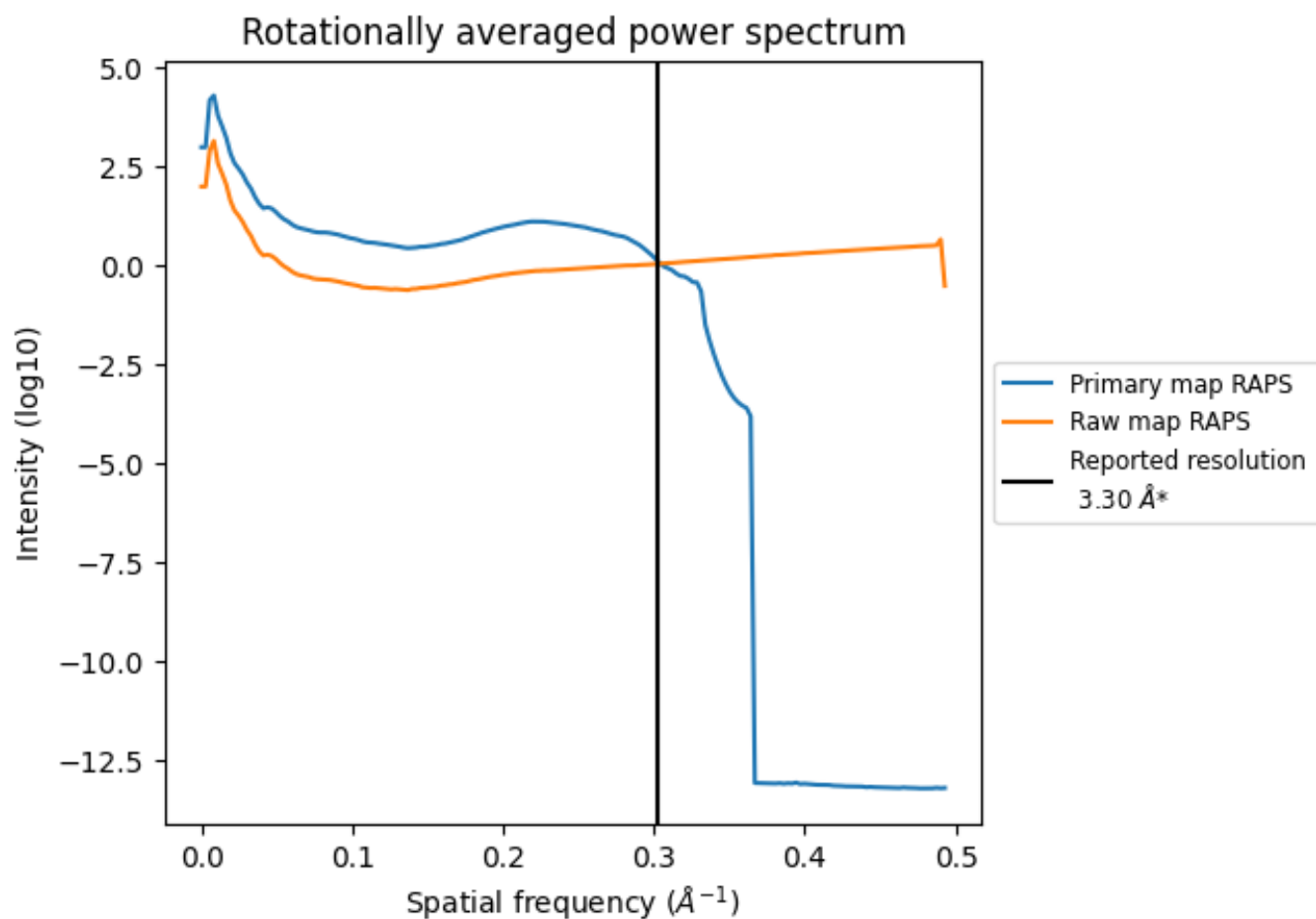
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 46 nm^3 ; this corresponds to an approximate mass of 41 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](#)

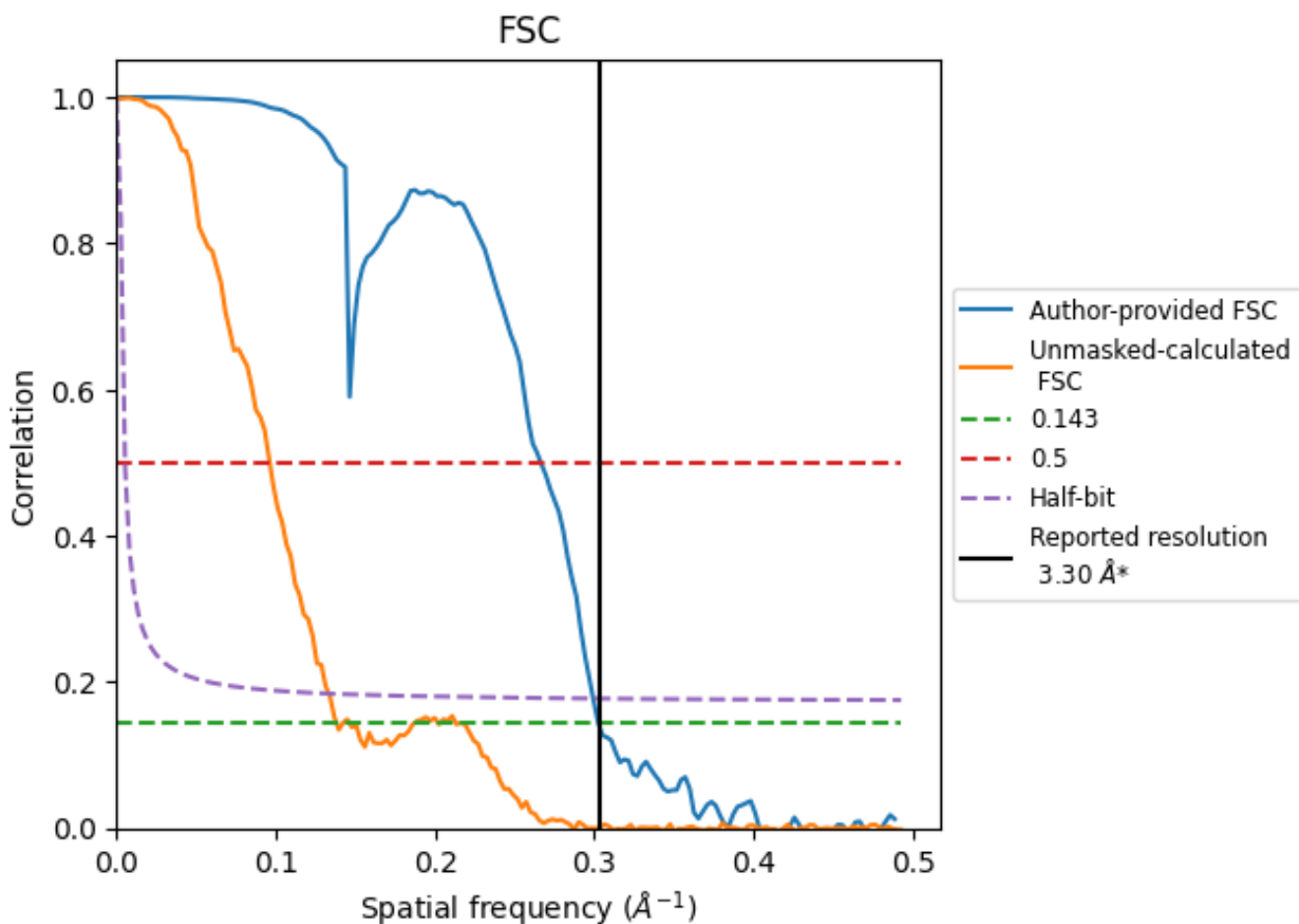


*Reported resolution corresponds to spatial frequency of 0.303 \AA^{-1}

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 0.303 Å⁻¹

8.2 Resolution estimates [i](#)

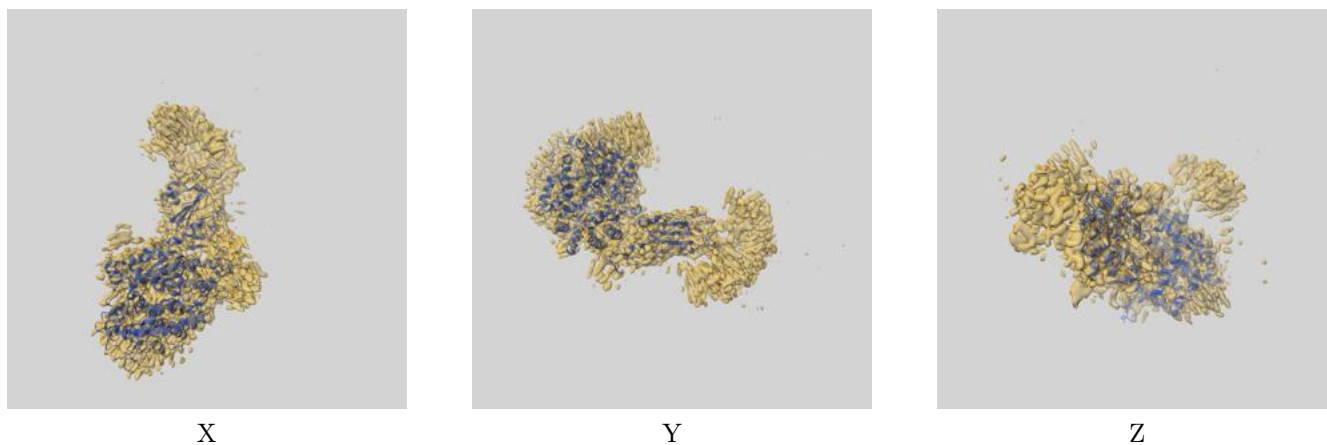
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	3.30	3.75	3.34
Unmasked-calculated*	7.23	10.38	7.50

*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 7.23 differs from the reported value 3.3 by more than 10 %

9 Map-model fit [i](#)

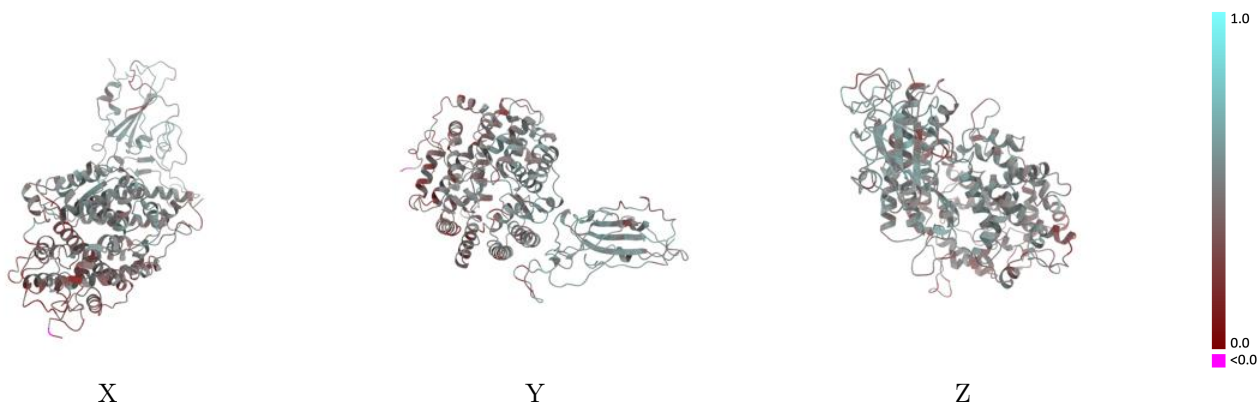
This section contains information regarding the fit between EMDB map EMD-15592 and PDB model 8AQW. Per-residue inclusion information can be found in section 3 on page 8.

9.1 Map-model overlay [i](#)



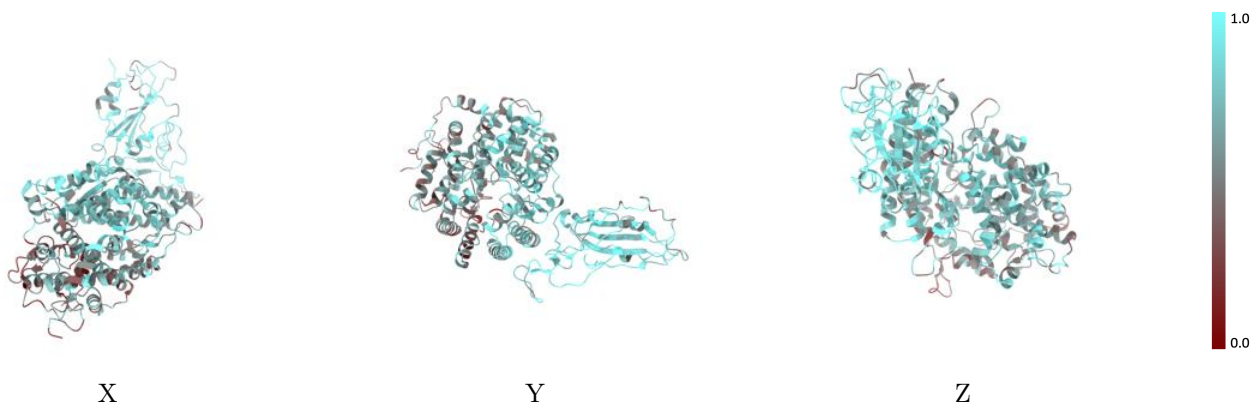
The images above show the 3D surface view of the map at the recommended contour level 0.172 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



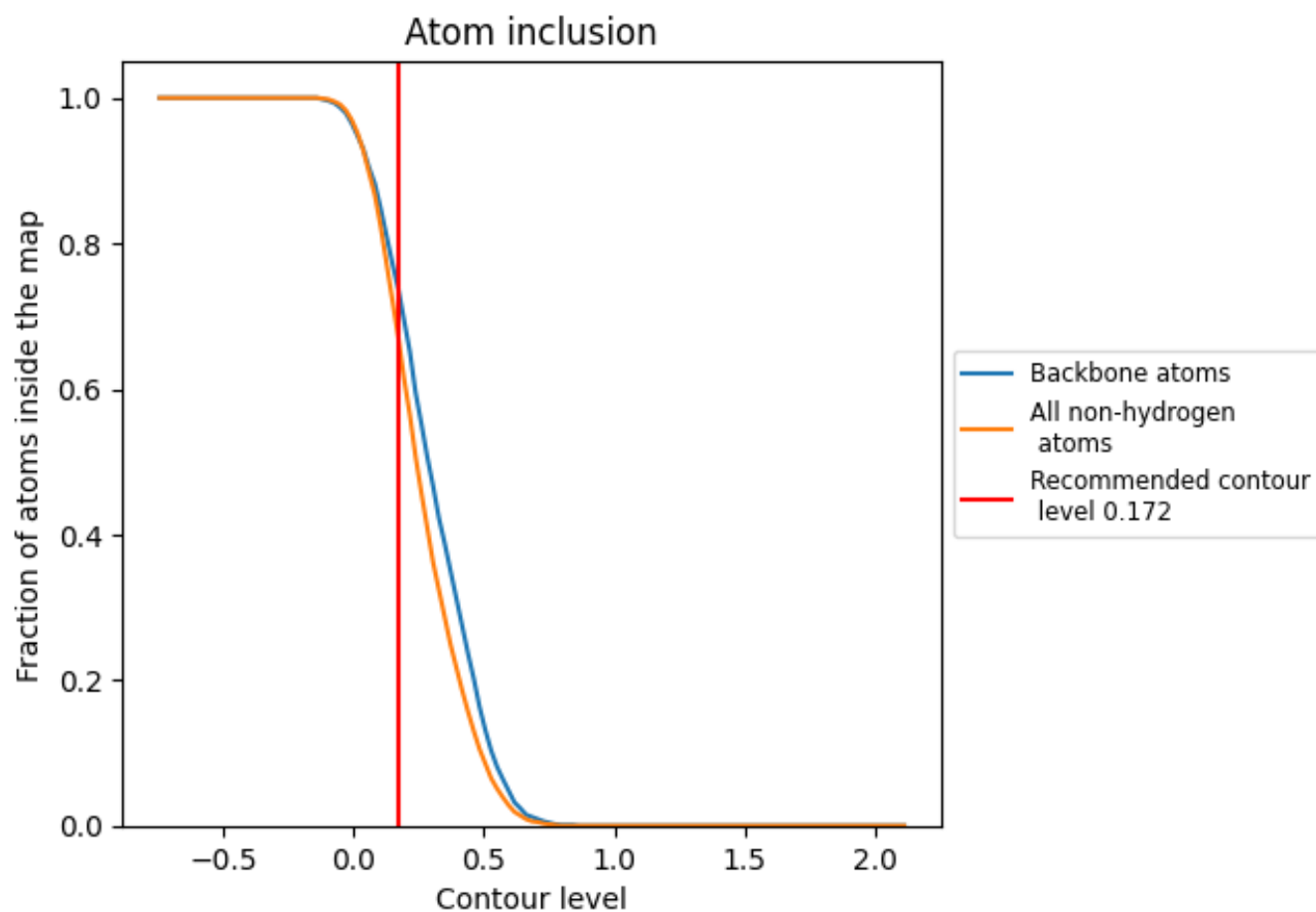
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.172).









9.4 Atom inclusion [i](#)



At the recommended contour level, 74% of all backbone atoms, 67% 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 (0.172) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6690	 0.4520
A	 0.6290	 0.4350
B	 0.7990	 0.5080
C	 0.5000	 0.3640

