



Full wwPDB EM Validation Report ⓘ

Oct 6, 2024 – 10:59 PM EDT

PDB ID : 8VGT
EMDB ID : EMD-43224
Title : Structure of the HKU1 RBD bound to the human TMPRSS2 receptor
Authors : Park, Y.J.; Seattle Structural Genomics Center for Infectious Disease (SSG-CID); Veesler, D.
Deposited on : 2023-12-27
Resolution : 2.90 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.dev113
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

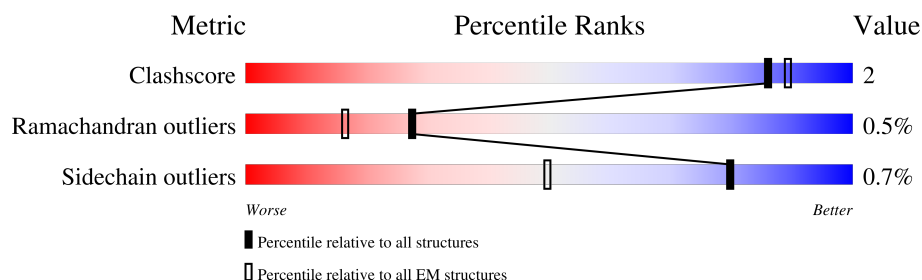
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.90 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	338	<div> <div>14%</div> <div>79%</div> <div>17%</div> </div>
2	B	416	<div> <div>6%</div> <div>79%</div> <div>16%</div> </div>
3	C	2	<div> <div>50%</div> <div>100%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4714 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 protein S1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	280	Total	C	N	O	S	0	0
			2040	1303	357	358	22		

There are 43 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	288	MET	-	initiating methionine	UNP Q5MQD0
A	289	GLY	-	expression tag	UNP Q5MQD0
A	290	ILE	-	expression tag	UNP Q5MQD0
A	291	LEU	-	expression tag	UNP Q5MQD0
A	292	PRO	-	expression tag	UNP Q5MQD0
A	293	SER	-	expression tag	UNP Q5MQD0
A	294	PRO	-	expression tag	UNP Q5MQD0
A	295	GLY	-	expression tag	UNP Q5MQD0
A	296	MET	-	expression tag	UNP Q5MQD0
A	297	PRO	-	expression tag	UNP Q5MQD0
A	298	ALA	-	expression tag	UNP Q5MQD0
A	299	LEU	-	expression tag	UNP Q5MQD0
A	300	LEU	-	expression tag	UNP Q5MQD0
A	301	SER	-	expression tag	UNP Q5MQD0
A	302	LEU	-	expression tag	UNP Q5MQD0
A	303	VAL	-	expression tag	UNP Q5MQD0
A	304	SER	-	expression tag	UNP Q5MQD0
A	305	LEU	-	expression tag	UNP Q5MQD0
A	306	LEU	-	expression tag	UNP Q5MQD0
A	307	SER	-	expression tag	UNP Q5MQD0
A	308	VAL	-	expression tag	UNP Q5MQD0
A	309	LEU	-	expression tag	UNP Q5MQD0
A	310	LEU	-	expression tag	UNP Q5MQD0
A	311	MET	-	expression tag	UNP Q5MQD0
A	312	GLY	-	expression tag	UNP Q5MQD0
A	313	CYS	-	expression tag	UNP Q5MQD0
A	314	VAL	-	expression tag	UNP Q5MQD0
A	315	ALA	-	expression tag	UNP Q5MQD0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	316	GLU	-	expression tag	UNP Q5MQD0
A	317	THR	-	expression tag	UNP Q5MQD0
A	318	GLY	-	expression tag	UNP Q5MQD0
A	319	THR	-	expression tag	UNP Q5MQD0
A	615	GLY	-	expression tag	UNP Q5MQD0
A	616	GLY	-	expression tag	UNP Q5MQD0
A	617	SER	-	expression tag	UNP Q5MQD0
A	618	HIS	-	expression tag	UNP Q5MQD0
A	619	HIS	-	expression tag	UNP Q5MQD0
A	620	HIS	-	expression tag	UNP Q5MQD0
A	621	HIS	-	expression tag	UNP Q5MQD0
A	622	HIS	-	expression tag	UNP Q5MQD0
A	623	HIS	-	expression tag	UNP Q5MQD0
A	624	HIS	-	expression tag	UNP Q5MQD0
A	625	HIS	-	expression tag	UNP Q5MQD0

- Molecule 2 is a protein called Transmembrane protease serine 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	349	Total	C	N	O	S	0	0
			2604	1671	461	447	25		

There are 41 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	91	MET	-	initiating methionine	UNP O15393
B	92	LYS	-	expression tag	UNP O15393
B	93	TRP	-	expression tag	UNP O15393
B	94	VAL	-	expression tag	UNP O15393
B	95	THR	-	expression tag	UNP O15393
B	96	PHE	-	expression tag	UNP O15393
B	97	ILE	-	expression tag	UNP O15393
B	98	SER	-	expression tag	UNP O15393
B	99	LEU	-	expression tag	UNP O15393
B	100	LEU	-	expression tag	UNP O15393
B	101	PHE	-	expression tag	UNP O15393
B	102	LEU	-	expression tag	UNP O15393
B	103	PHE	-	expression tag	UNP O15393
B	104	SER	-	expression tag	UNP O15393
B	105	SER	-	expression tag	UNP O15393
B	106	ALA	-	expression tag	UNP O15393
B	107	TYR	-	expression tag	UNP O15393

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Chain	Residue	Modelled	Actual	Comment	Reference
B	108	SER	-	expression tag	UNP O15393
B	?	-	SER	deletion	UNP O15393
B	251	ASP	SER	engineered mutation	UNP O15393
B	252	ASP	ARG	engineered mutation	UNP O15393
B	253	ASP	GLN	engineered mutation	UNP O15393
B	254	ASP	SER	engineered mutation	UNP O15393
B	255	LYS	ARG	engineered mutation	UNP O15393
B	441	ALA	SER	engineered mutation	UNP O15393
B	447	CYS	THR	engineered mutation	UNP O15393
B	493	ASP	-	expression tag	UNP O15393
B	494	ASP	-	expression tag	UNP O15393
B	495	ASP	-	expression tag	UNP O15393
B	496	ASP	-	expression tag	UNP O15393
B	497	LYS	-	expression tag	UNP O15393
B	498	SER	-	expression tag	UNP O15393
B	499	GLY	-	expression tag	UNP O15393
B	500	HIS	-	expression tag	UNP O15393
B	501	HIS	-	expression tag	UNP O15393
B	502	HIS	-	expression tag	UNP O15393
B	503	HIS	-	expression tag	UNP O15393
B	504	HIS	-	expression tag	UNP O15393
B	505	HIS	-	expression tag	UNP O15393
B	506	HIS	-	expression tag	UNP O15393
B	507	HIS	-	expression tag	UNP O15393

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

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

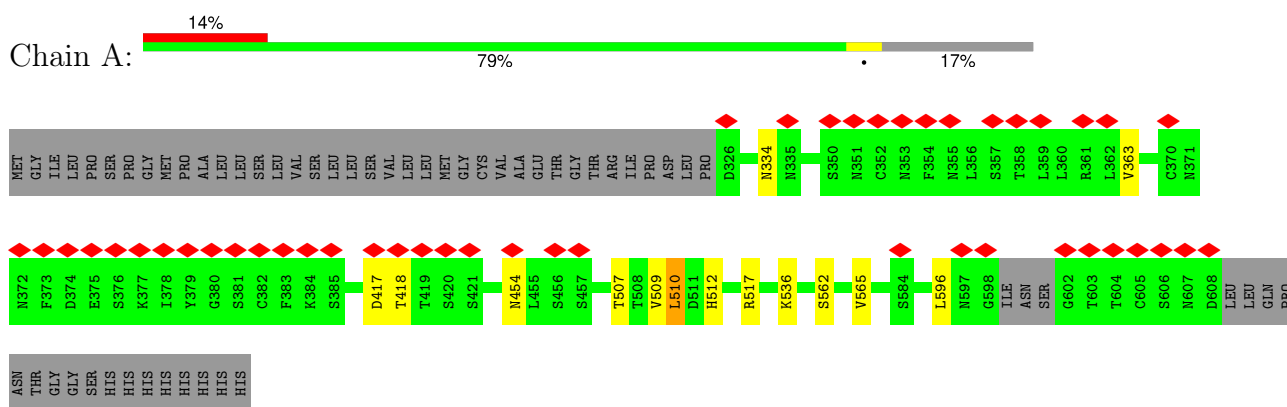


Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	

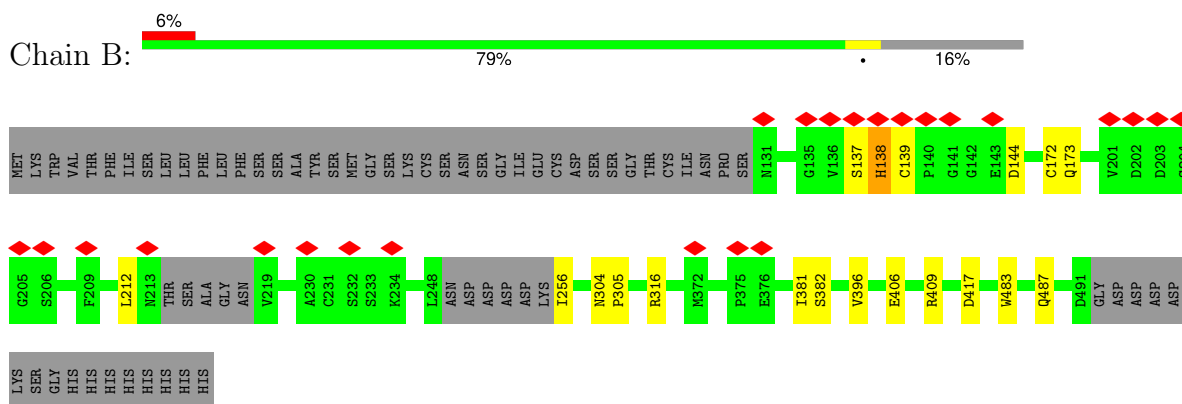
3 Residue-property plots

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.

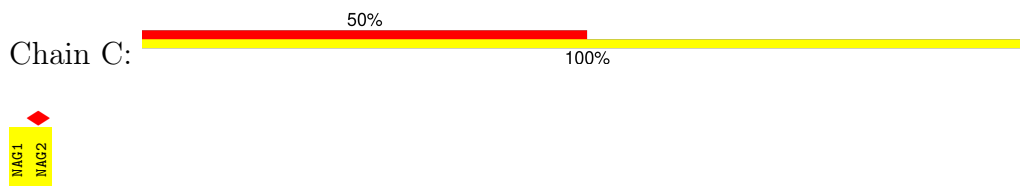
- Molecule 1: Spike protein S1



- Molecule 2: Transmembrane protease serine 2



- Molecule 3: 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	810357	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	10.827	Depositor
Minimum map value	-6.958	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.126	Depositor
Recommended contour level	1.1	Depositor
Map size (Å)	280.0, 280.0, 280.0	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.0, 1.0, 1.0	Depositor

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.57	0/2095	0.58	0/2865
2	B	0.60	0/2676	0.57	0/3648
All	All	0.58	0/4771	0.57	0/6513

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2040	0	1718	8	0
2	B	2604	0	2387	10	0
3	C	28	0	25	0	0
4	A	42	0	39	0	0
All	All	4714	0	4169	18	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 2.

All (18) 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:B:316:ARG:HG2	2:B:396:VAL:HG12	1.51	0.90
1:A:334:ASN:OD1	1:A:363:VAL:HG12	1.88	0.73
1:A:507:THR:HG21	1:A:517:ARG:HG3	1.72	0.72
1:A:596:LEU:N	1:A:596:LEU:HD22	2.15	0.61
2:B:137:SER:O	2:B:139:CYS:N	2.37	0.57
2:B:256:ILE:HD12	2:B:381:ILE:HD12	1.88	0.56
1:A:509:VAL:O	1:A:512:HIS:HD2	1.96	0.48
2:B:256:ILE:HG13	2:B:382:SER:HA	1.95	0.48
1:A:562:SER:O	1:A:565:VAL:HG22	2.14	0.47
1:A:509:VAL:O	1:A:510:LEU:HB2	2.15	0.47
2:B:137:SER:O	2:B:138:HIS:C	2.55	0.45
2:B:483:TRP:O	2:B:487:GLN:NE2	2.49	0.45
1:A:417:ASP:OD2	1:A:418:THR:N	2.51	0.43
2:B:406:GLU:HG3	2:B:409:ARG:H	1.85	0.42
1:A:596:LEU:N	1:A:596:LEU:CD2	2.80	0.42
2:B:304:ASN:HA	2:B:305:PRO:HD3	1.87	0.41
2:B:144:ASP:N	2:B:144:ASP:OD1	2.52	0.41
2:B:172:CYS:SG	2:B:173:GLN:N	2.94	0.41

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	276/338 (82%)	265 (96%)	10 (4%)	1 (0%)	30	60
2	B	343/416 (82%)	327 (95%)	14 (4%)	2 (1%)	22	52
All	All	619/754 (82%)	592 (96%)	24 (4%)	3 (0%)	27	56

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	510	LEU

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Mol	Chain	Res	Type
2	B	138	HIS
2	B	417	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	186/308 (60%)	184 (99%)	2 (1%)	70	90
2	B	244/355 (69%)	243 (100%)	1 (0%)	89	97
All	All	430/663 (65%)	427 (99%)	3 (1%)	80	94

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

Mol	Chain	Res	Type
1	A	454	ASN
1	A	536	LYS
2	B	212	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	512	HIS
2	B	307	HIS
2	B	334	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

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	1.19	1 (7%)	17,19,21	1.12	1 (5%)
3	NAG	C	2	3	14,14,15	1.13	2 (14%)	17,19,21	0.95	1 (5%)

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	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	NAG	C1-C2	3.27	1.56	1.52
3	C	2	NAG	C1-C2	3.00	1.56	1.52
3	C	2	NAG	O5-C5	2.23	1.47	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	NAG	C8-C7-N2	2.66	120.54	116.12
3	C	2	NAG	C8-C7-N2	2.30	119.92	116.12

There are no chirality outliers.

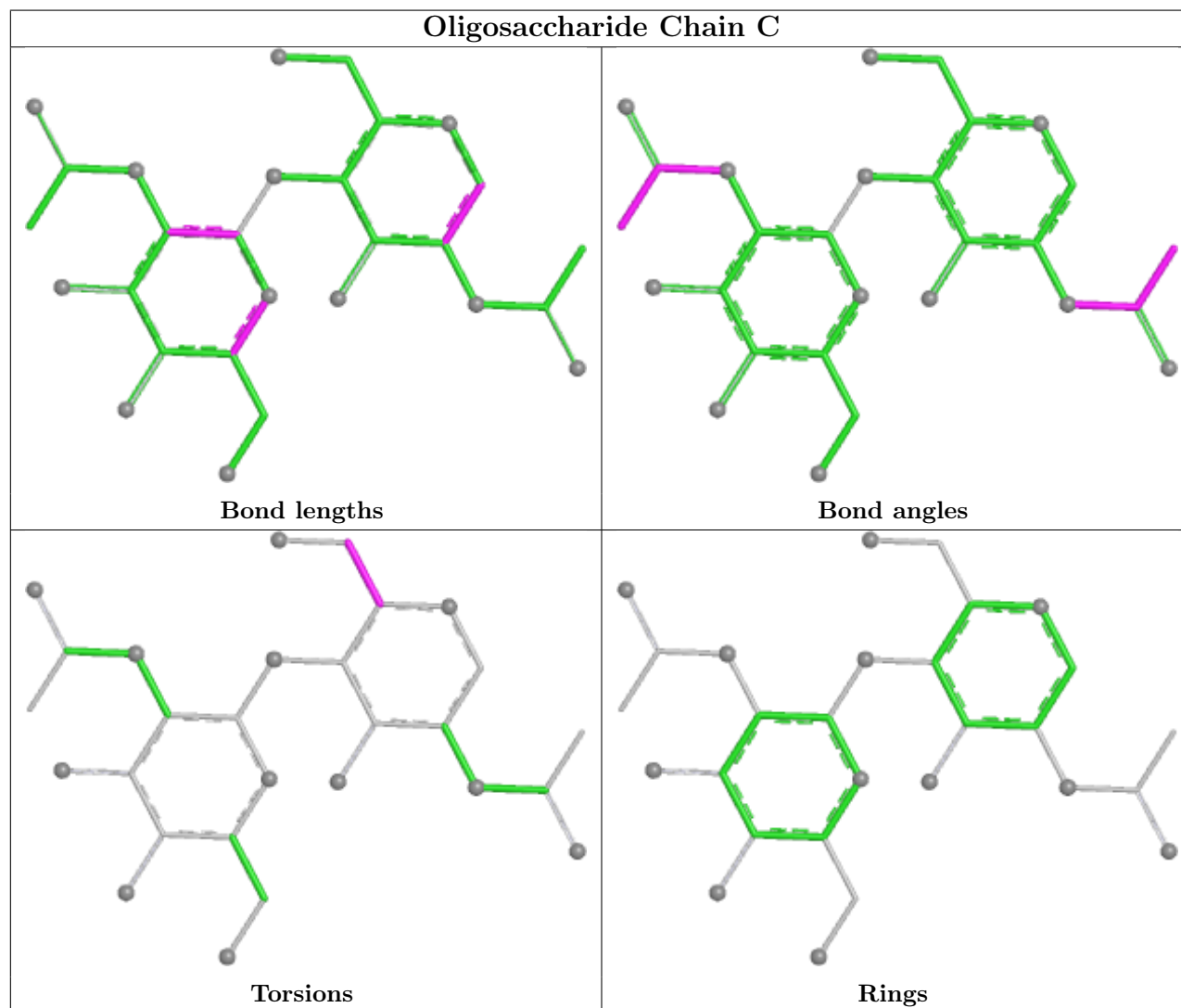
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1	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 ⓘ

3 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
4	NAG	A	702	1	14,14,15	1.19	1 (7%)	17,19,21	1.04	1 (5%)
4	NAG	A	703	1	14,14,15	1.14	1 (7%)	17,19,21	1.01	1 (5%)
4	NAG	A	701	1	14,14,15	1.17	1 (7%)	17,19,21	1.05	1 (5%)

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	A	702	1	-	0/6/23/26	0/1/1/1
4	NAG	A	703	1	-	2/6/23/26	0/1/1/1
4	NAG	A	701	1	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	702	NAG	C1-C2	3.61	1.57	1.52
4	A	703	NAG	C1-C2	3.45	1.57	1.52
4	A	701	NAG	C1-C2	3.38	1.57	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	701	NAG	C8-C7-N2	2.57	120.37	116.12
4	A	702	NAG	C8-C7-N2	2.43	120.15	116.12
4	A	703	NAG	C8-C7-N2	2.25	119.84	116.12

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	703	NAG	O5-C5-C6-O6
4	A	703	NAG	C4-C5-C6-O6

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.

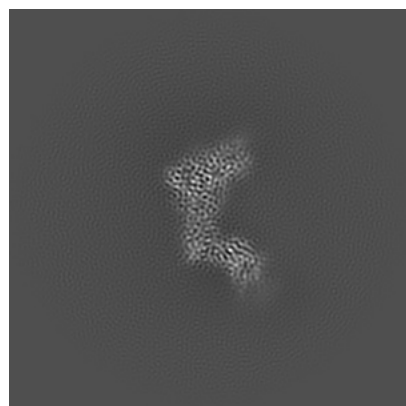
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-43224. These allow visual inspection of the internal detail of the map and identification of artifacts.

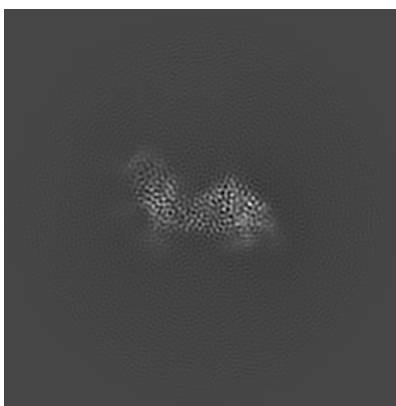
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

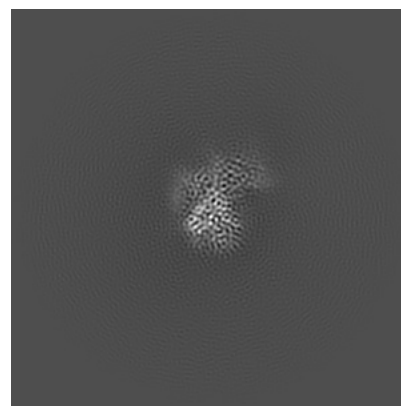
6.1.1 Primary map



X

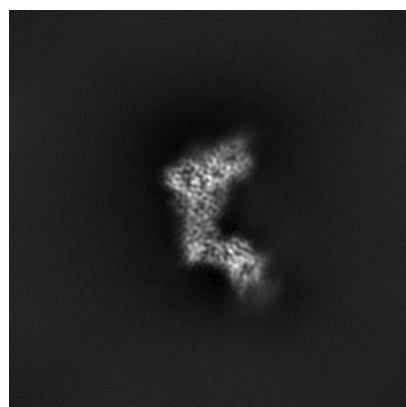


Y

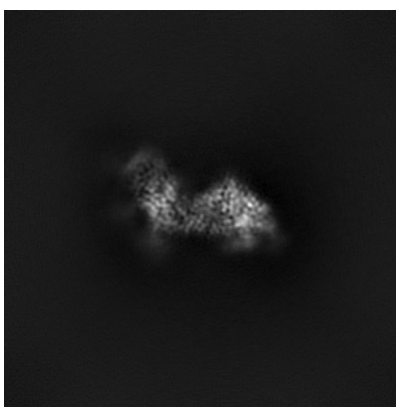


Z

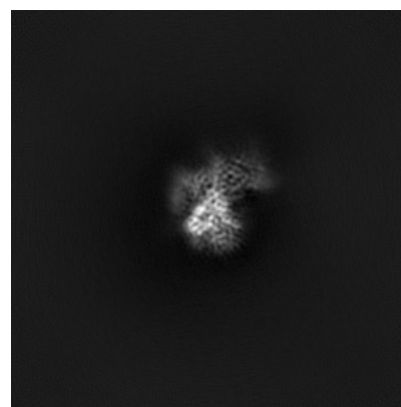
6.1.2 Raw map



X



Y

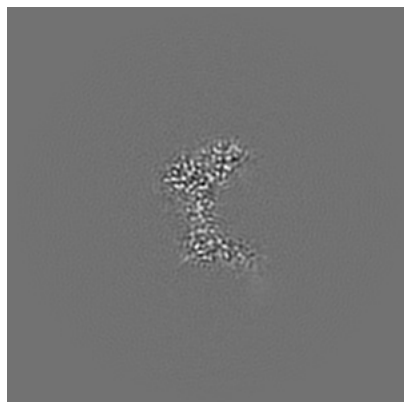


Z

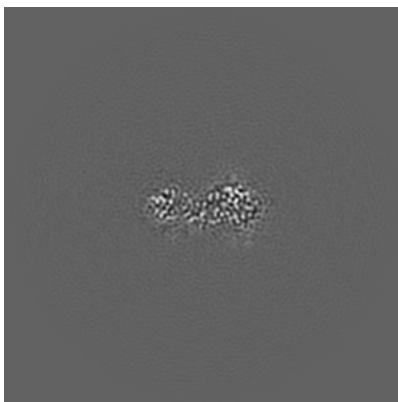
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

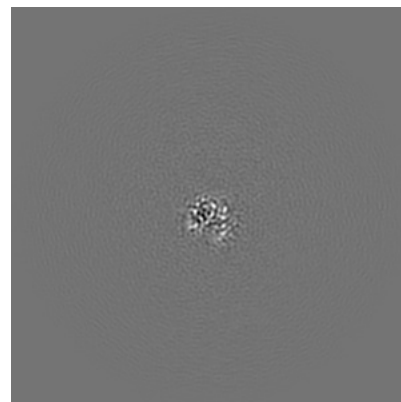
6.2.1 Primary map



X Index: 140

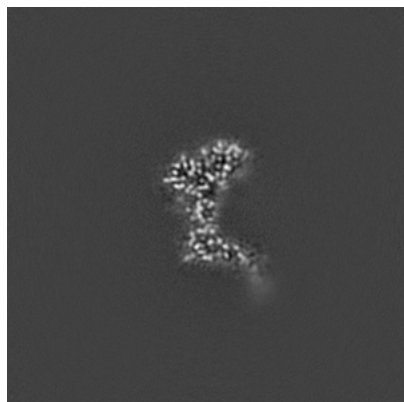


Y Index: 140

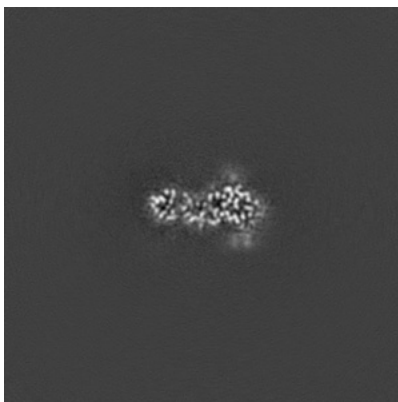


Z Index: 140

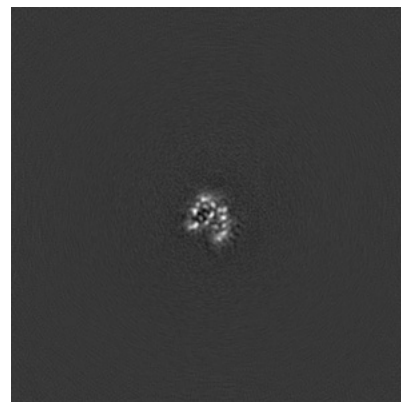
6.2.2 Raw map



X Index: 140



Y Index: 140

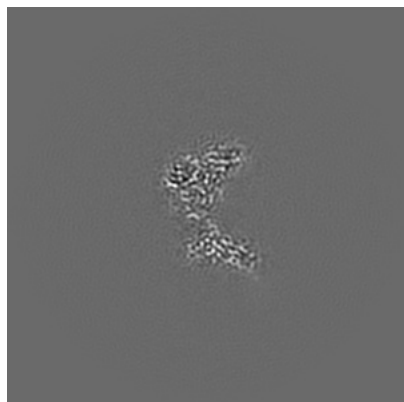


Z Index: 140

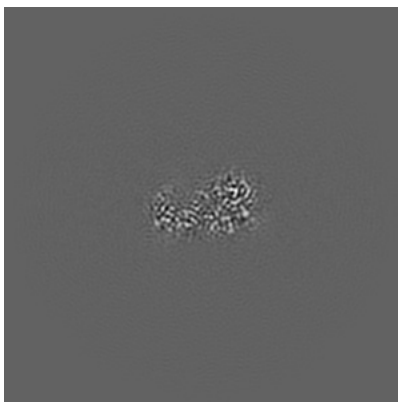
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

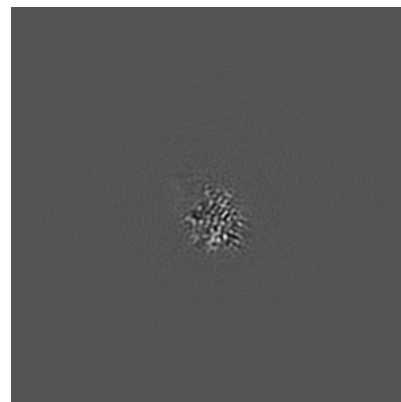
6.3.1 Primary map



X Index: 143

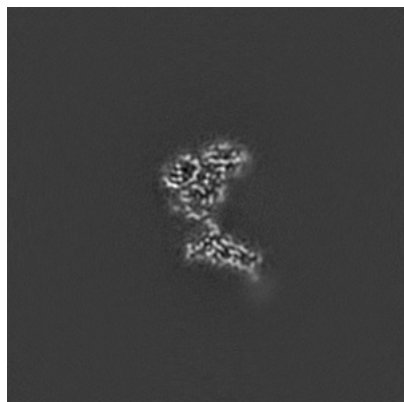


Y Index: 131

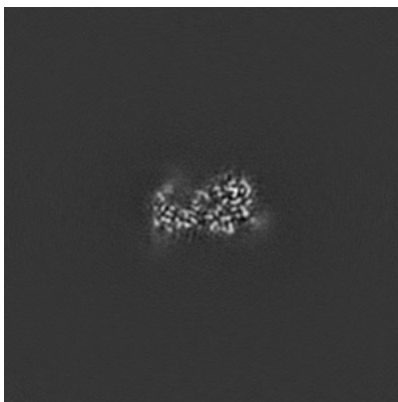


Z Index: 159

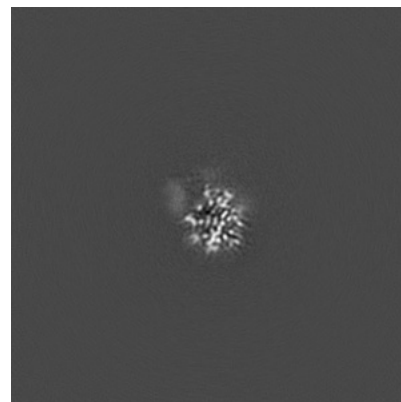
6.3.2 Raw map



X Index: 143



Y Index: 131

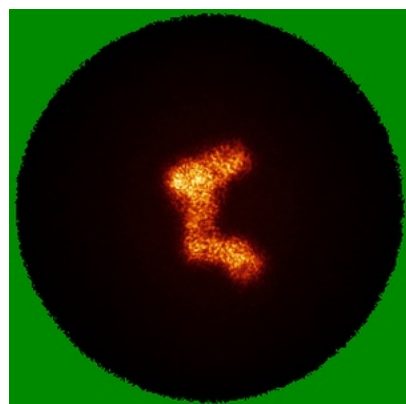


Z Index: 159

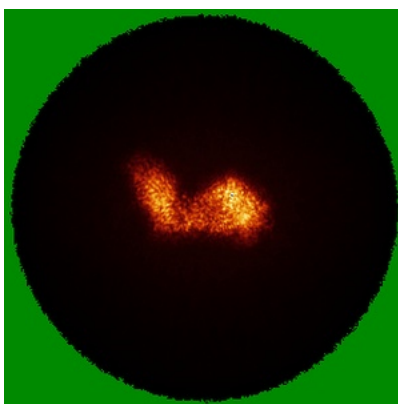
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) ⓘ

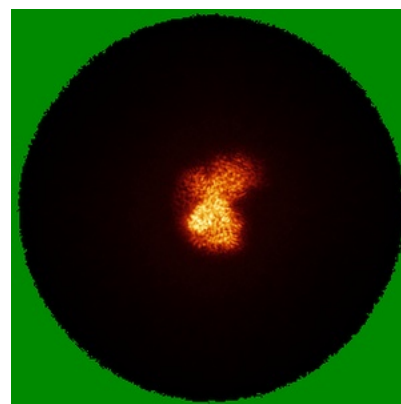
6.4.1 Primary map



X

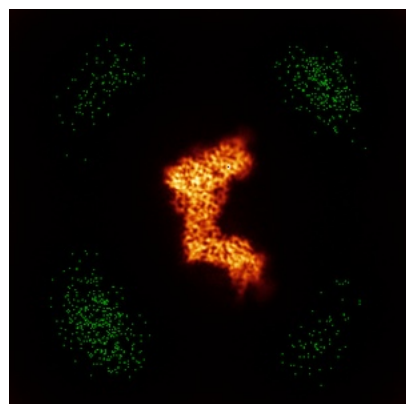


Y

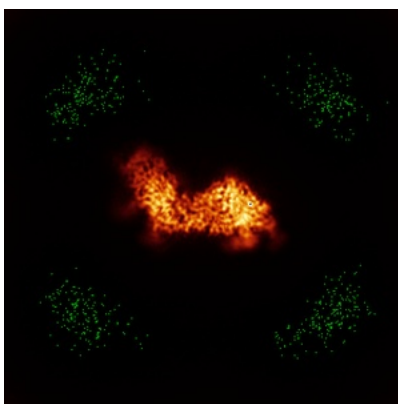


Z

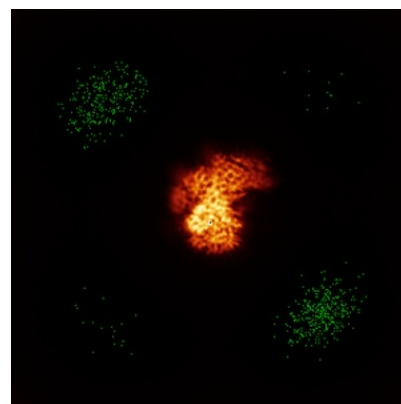
6.4.2 Raw map



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.

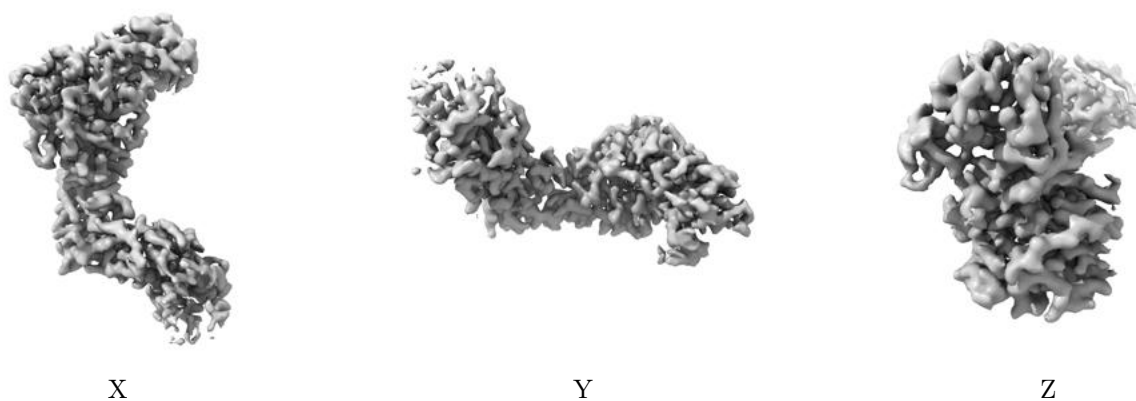
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 1.1. 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.

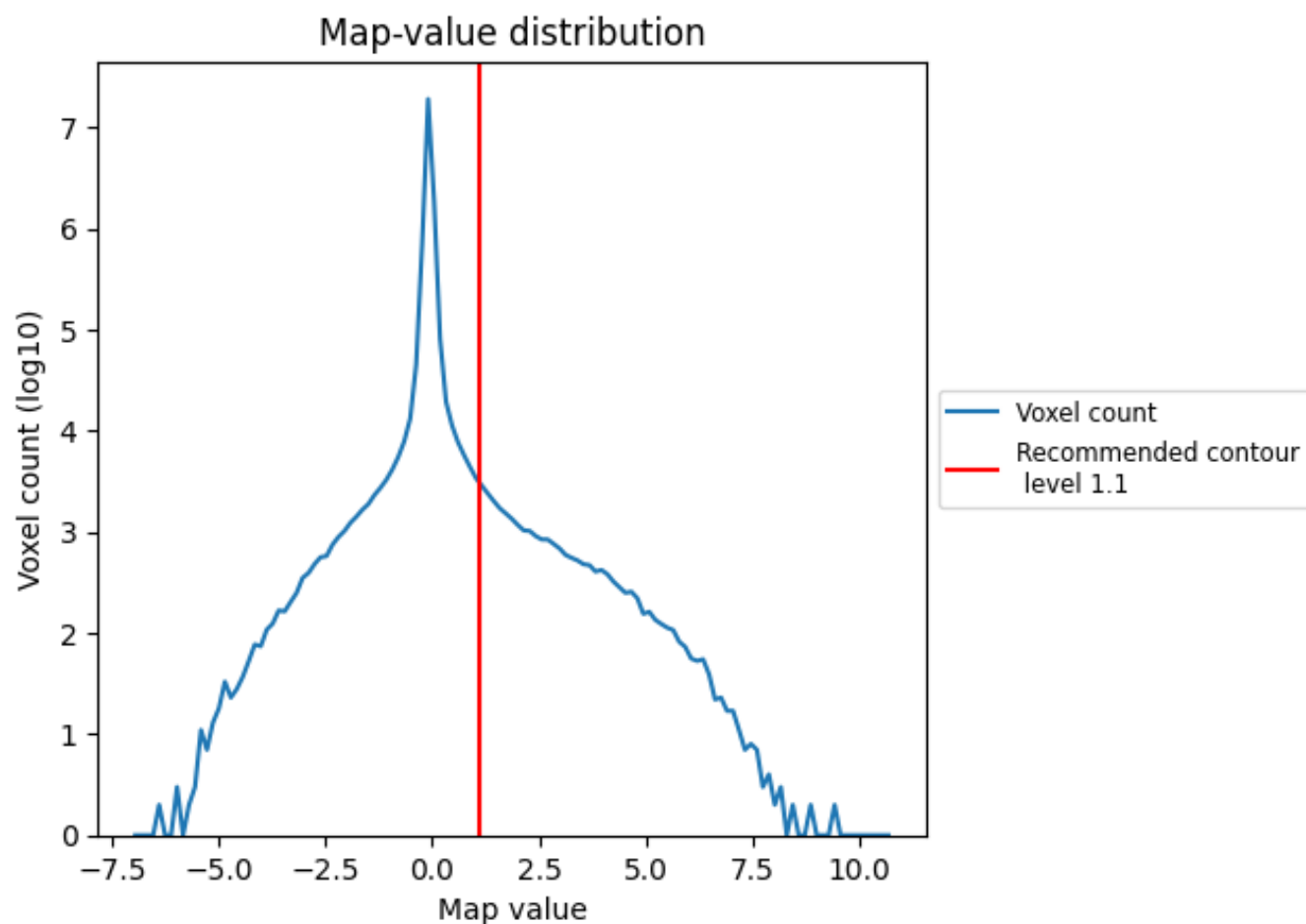
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

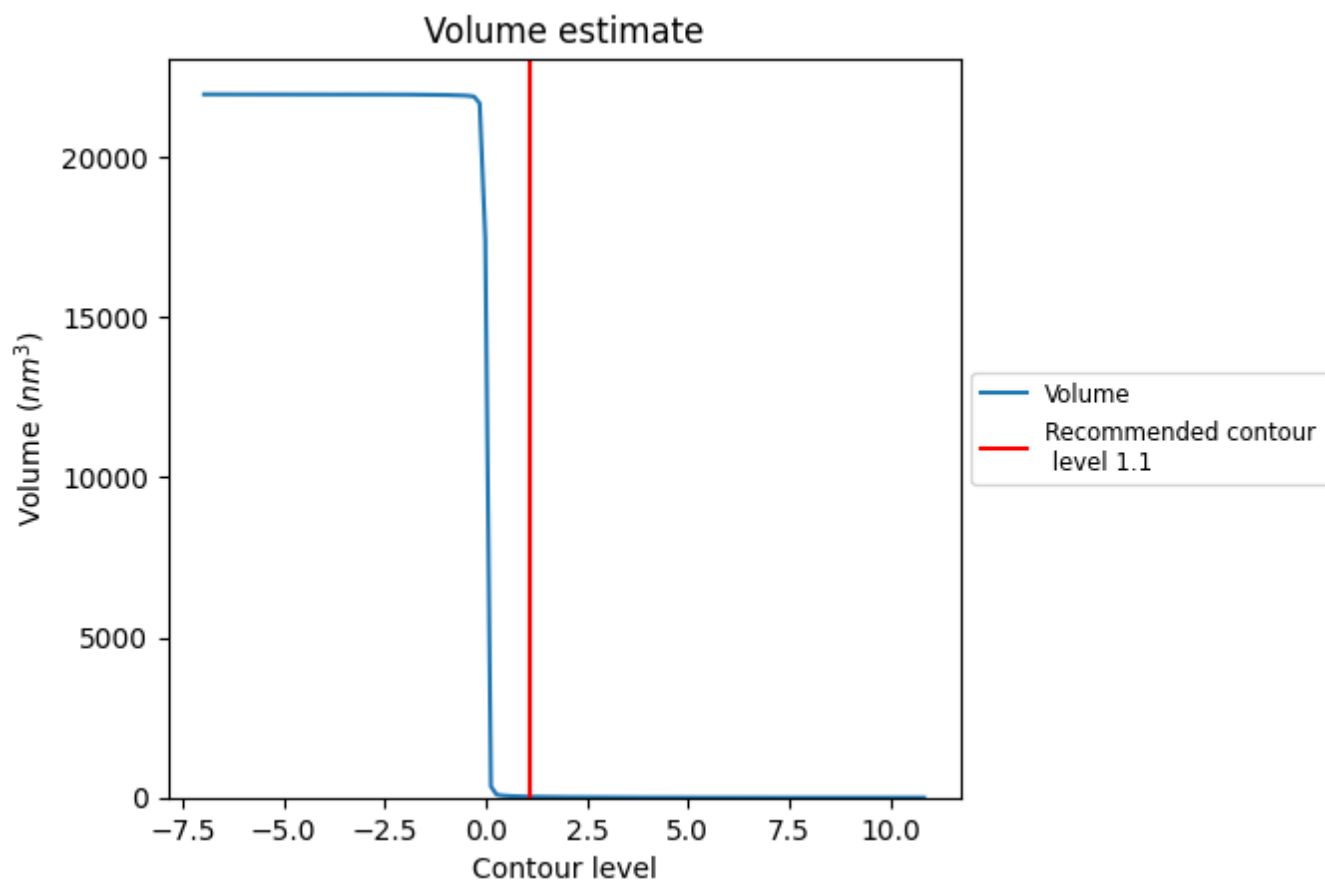
This section contains the results of statistical analysis of the map.

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.

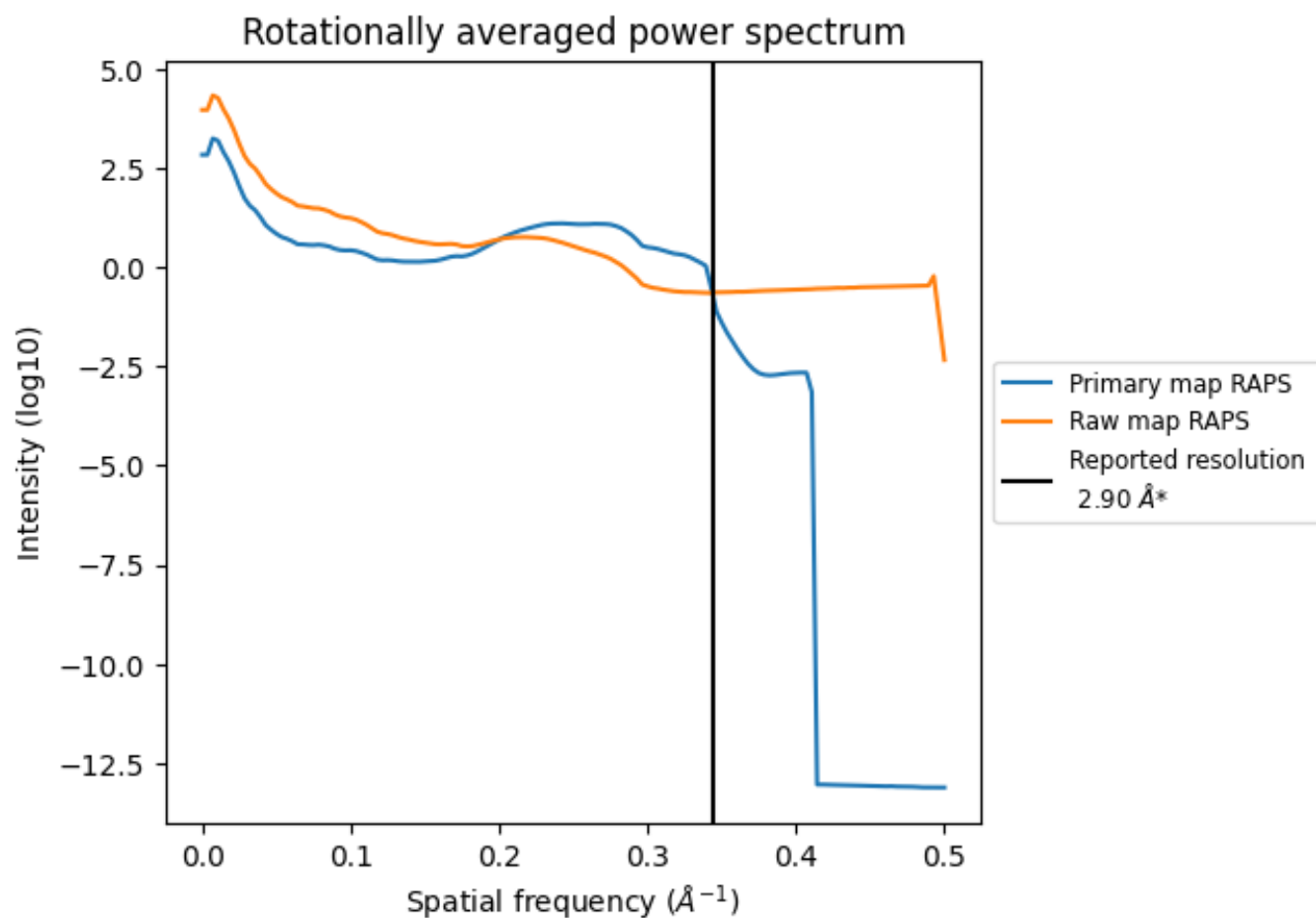
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 27 nm³; this corresponds to an approximate mass of 24 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

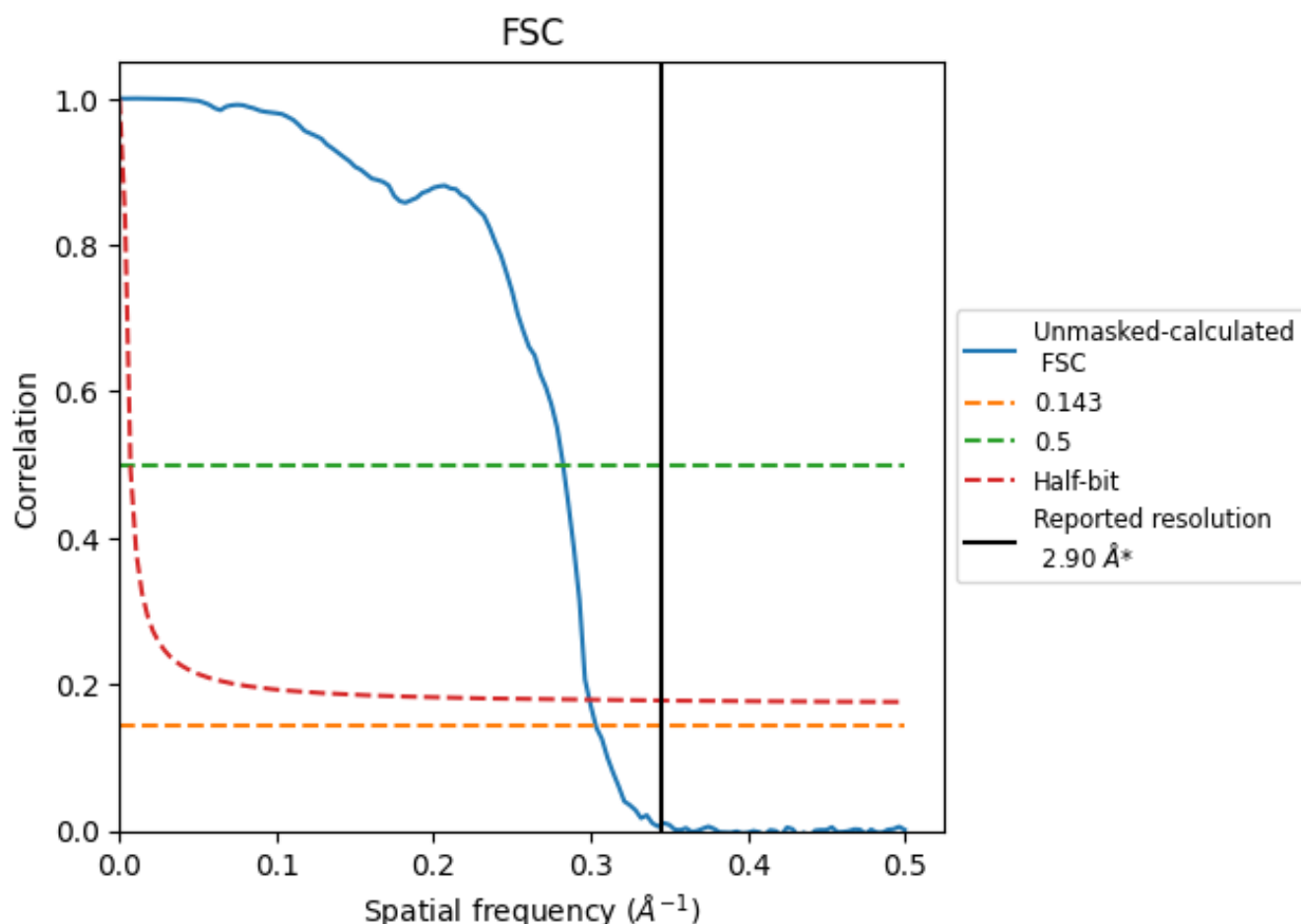


*Reported resolution corresponds to spatial frequency of 0.345 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.345 \AA^{-1}

8.2 Resolution estimates [i](#)

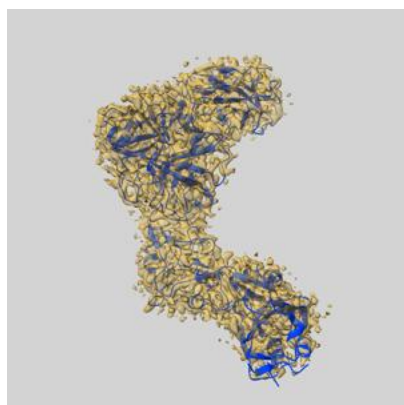
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.90	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.30	3.54	3.34

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.30 differs from the reported value 2.9 by more than 10 %

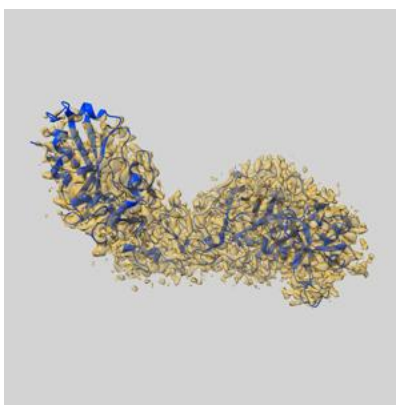
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-43224 and PDB model 8VGT. Per-residue inclusion information can be found in [section 3](#) on [page 7](#).

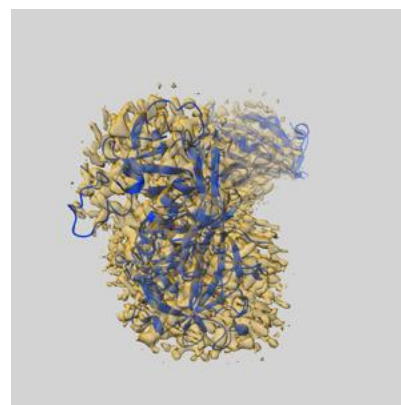
9.1 Map-model overlay [i](#)



X



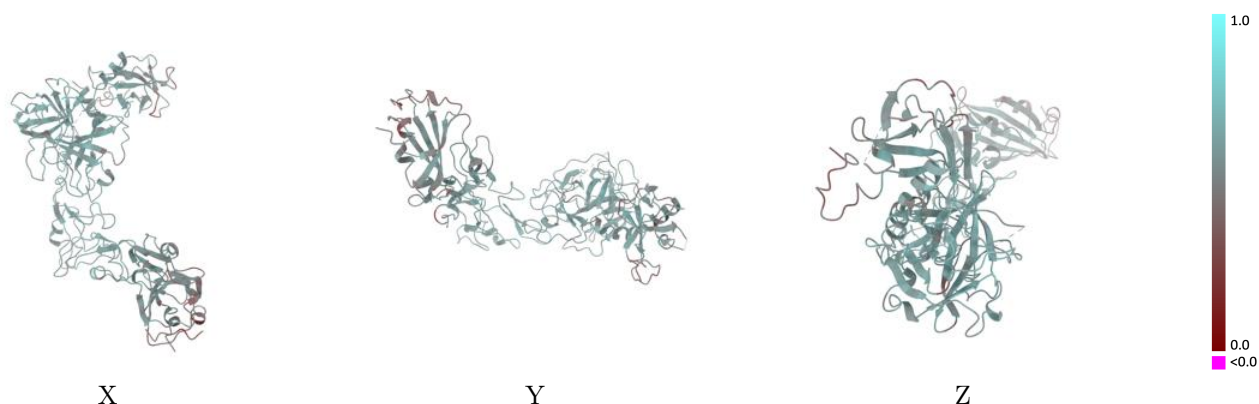
Y



Z

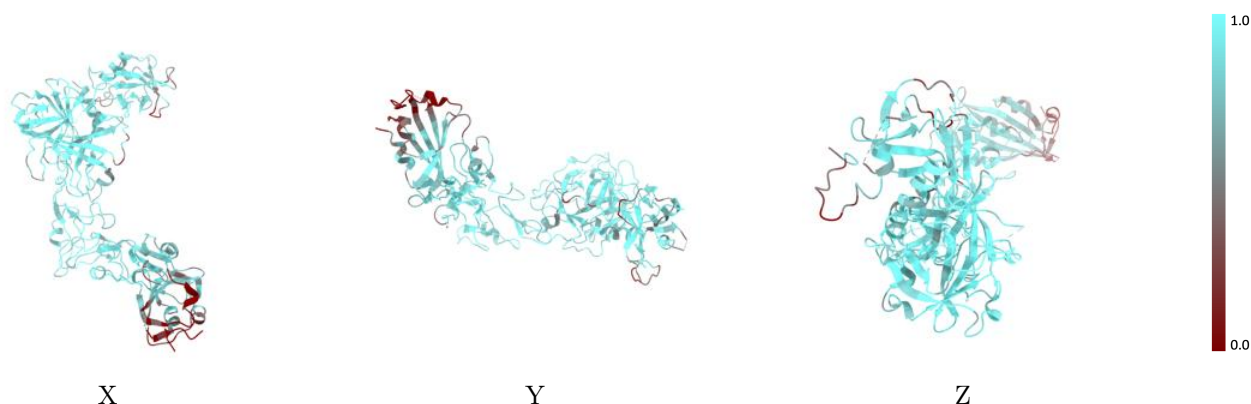
The images above show the 3D surface view of the map at the recommended contour level 1.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



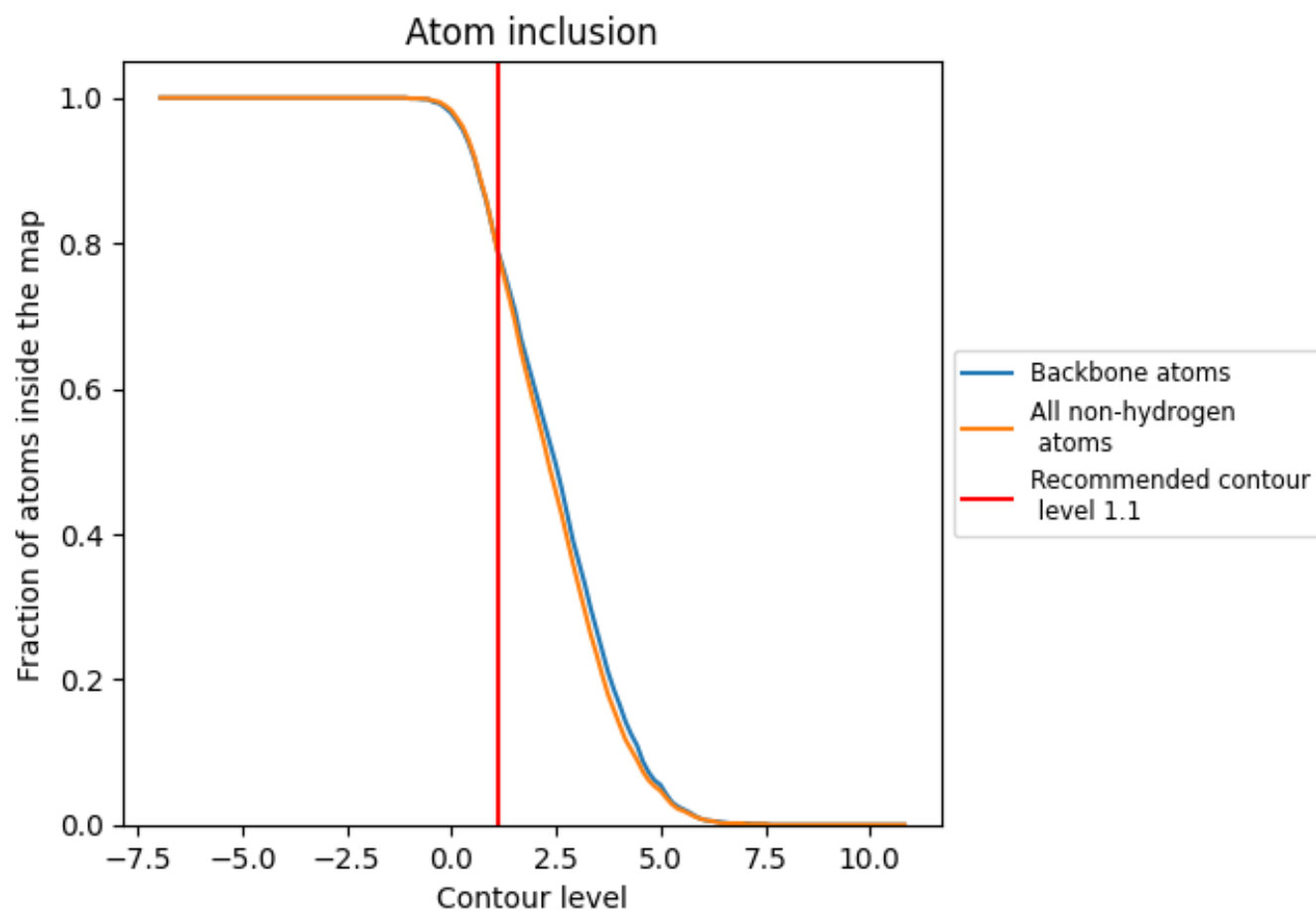
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (1.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 79% 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 (1.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.7880	<div></div> 0.5530
A	<div></div> 0.7260	<div></div> 0.5390
B	<div></div> 0.8440	<div></div> 0.5670
C	<div></div> 0.2500	<div></div> 0.2550

