



## Full wwPDB EM Validation Report ⓘ

Dec 7, 2022 – 12:34 PM JST

PDB ID : 7DZW  
EMDB ID : EMD-30915  
Title : Apo spike protein from SARS-CoV2  
Authors : Liu, Y.; Soh, W.T.; Li, S.; Kishikawa, J.; Hirose, M.; Kato, T.; Standley, D.;  
Okada, M.; Arase, H.  
Deposited on : 2021-01-26  
Resolution : 3.45 Å(reported)  
Based on initial model : 7KEB

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

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

A user guide is available at

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

The types of validation reports are described at

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

---

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

EMDB validation analysis : 0.0.1.dev43  
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) : 2.31.3

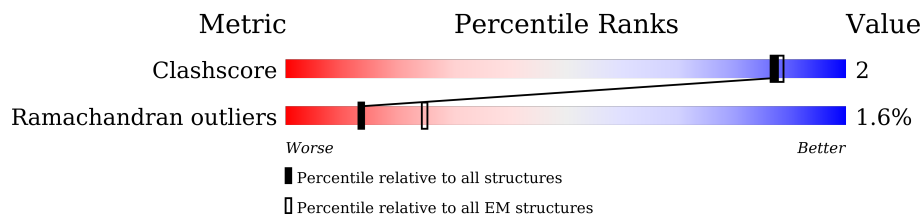
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.45 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

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

Mol	Chain	Length	Quality of chain
1	A	1249	<div> <div>18%</div> <div>87%</div> <div>10%</div> </div>
1	B	1249	<div> <div>25%</div> <div>86%</div> <div>10%</div> </div>
1	C	1249	<div> <div>19%</div> <div>88%</div> <div>10%</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 13452 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.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	1121	Total	C	N	O	0	0
			4484	2242	1121	1121		
1	B	1121	Total	C	N	O	0	0
			4484	2242	1121	1121		
1	C	1121	Total	C	N	O	0	0
			4484	2242	1121	1121		

There are 42 discrepancies between the modelled and reference sequences:

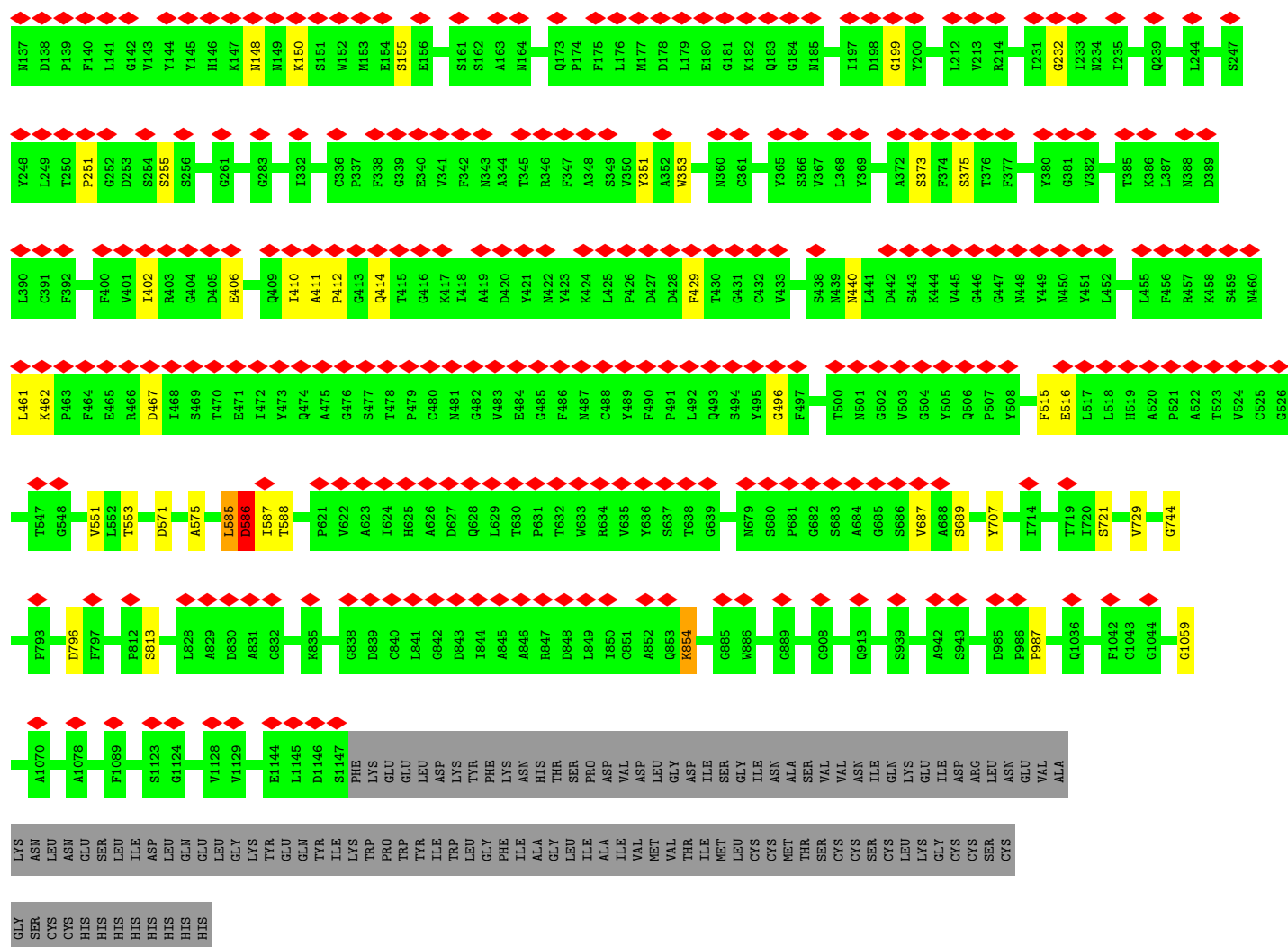
Chain	Residue	Modelled	Actual	Comment	Reference
A	614	GLY	ASP	engineered mutation	UNP P0DTC2
A	682	GLY	ARG	engineered mutation	UNP P0DTC2
A	683	SER	ARG	engineered mutation	UNP P0DTC2
A	685	GLY	ARG	engineered mutation	UNP P0DTC2
A	986	PRO	LYS	engineered mutation	UNP P0DTC2
A	987	PRO	VAL	engineered mutation	UNP P0DTC2
A	1255	HIS	-	expression tag	UNP P0DTC2
A	1256	HIS	-	expression tag	UNP P0DTC2
A	1257	HIS	-	expression tag	UNP P0DTC2
A	1258	HIS	-	expression tag	UNP P0DTC2
A	1259	HIS	-	expression tag	UNP P0DTC2
A	1260	HIS	-	expression tag	UNP P0DTC2
A	1261	HIS	-	expression tag	UNP P0DTC2
A	1262	HIS	-	expression tag	UNP P0DTC2
B	614	GLY	ASP	engineered mutation	UNP P0DTC2
B	682	GLY	ARG	engineered mutation	UNP P0DTC2
B	683	SER	ARG	engineered mutation	UNP P0DTC2
B	685	GLY	ARG	engineered mutation	UNP P0DTC2
B	986	PRO	LYS	engineered mutation	UNP P0DTC2
B	987	PRO	VAL	engineered mutation	UNP P0DTC2
B	1255	HIS	-	expression tag	UNP P0DTC2
B	1256	HIS	-	expression tag	UNP P0DTC2
B	1257	HIS	-	expression tag	UNP P0DTC2
B	1258	HIS	-	expression tag	UNP P0DTC2

*Continued on next page...*

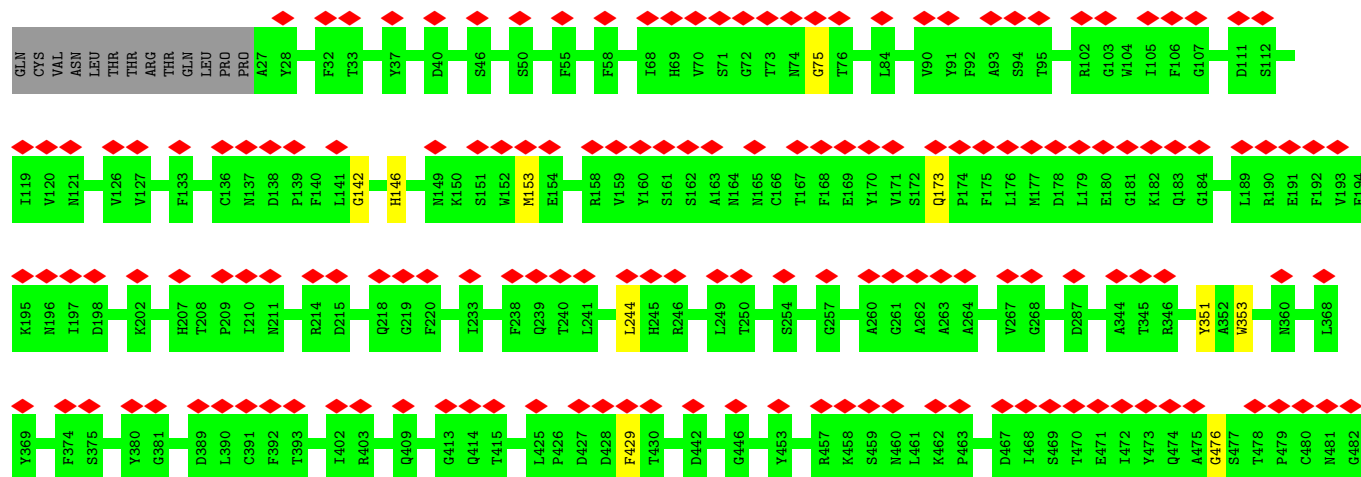
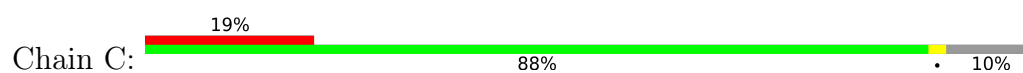
*Continued from previous page...*

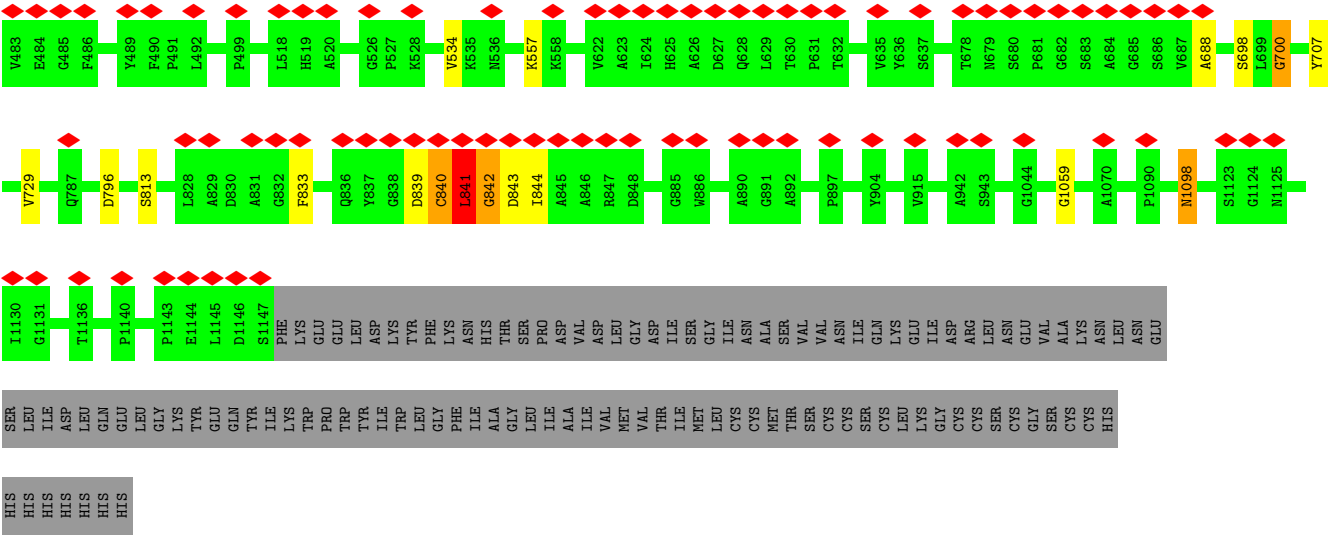
Chain	Residue	Modelled	Actual	Comment	Reference
B	1259	HIS	-	expression tag	UNP P0DTC2
B	1260	HIS	-	expression tag	UNP P0DTC2
B	1261	HIS	-	expression tag	UNP P0DTC2
B	1262	HIS	-	expression tag	UNP P0DTC2
C	614	GLY	ASP	engineered mutation	UNP P0DTC2
C	682	GLY	ARG	engineered mutation	UNP P0DTC2
C	683	SER	ARG	engineered mutation	UNP P0DTC2
C	685	GLY	ARG	engineered mutation	UNP P0DTC2
C	986	PRO	LYS	engineered mutation	UNP P0DTC2
C	987	PRO	VAL	engineered mutation	UNP P0DTC2
C	1255	HIS	-	expression tag	UNP P0DTC2
C	1256	HIS	-	expression tag	UNP P0DTC2
C	1257	HIS	-	expression tag	UNP P0DTC2
C	1258	HIS	-	expression tag	UNP P0DTC2
C	1259	HIS	-	expression tag	UNP P0DTC2
C	1260	HIS	-	expression tag	UNP P0DTC2
C	1261	HIS	-	expression tag	UNP P0DTC2
C	1262	HIS	-	expression tag	UNP P0DTC2





### • Molecule 1: Spike glycoprotein





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	120442	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)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.151	Depositor
Minimum map value	-0.503	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.039	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	316.8, 316.8, 316.8	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.88, 0.88, 0.88	Depositor



## 5 Model quality

### 5.1 Standard geometry

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	0.86	0/4483	0.92	6/5602 (0.1%)
1	B	0.88	0/4483	1.00	17/5602 (0.3%)
1	C	0.88	3/4483 (0.1%)	0.94	10/5602 (0.2%)
All	All	0.87	3/13449 (0.0%)	0.95	33/16806 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	842	GLY	CA-C	9.70	1.67	1.51
1	C	842	GLY	N-CA	6.71	1.56	1.46
1	C	841	LEU	C-N	5.29	1.42	1.33

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	842	GLY	C-N-CA	15.64	160.80	121.70
1	A	141	LEU	C-N-CA	-9.68	101.97	122.30
1	B	515	PHE	C-N-CA	8.64	143.31	121.70
1	B	854	LYS	N-CA-C	-8.49	88.07	111.00
1	C	840	CYS	C-N-CA	8.47	142.88	121.70
1	B	496	GLY	C-N-CA	7.70	140.95	121.70
1	B	461	LEU	C-N-CA	-7.62	102.66	121.70
1	B	39	PRO	N-CA-C	-6.48	95.26	112.10
1	B	586	ASP	O-C-N	-6.22	112.74	122.70
1	B	586	ASP	CA-C-N	6.06	130.54	117.20

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	794	ILE	C-N-CA	6.05	136.82	121.70
1	B	373	SER	C-N-CA	6.04	136.81	121.70
1	A	793	PRO	N-CA-C	-5.95	96.63	112.10
1	B	251	PRO	N-CA-C	-5.74	97.19	112.10
1	B	585	LEU	C-N-CA	5.69	135.93	121.70
1	C	842	GLY	N-CA-C	5.67	127.27	113.10
1	C	1098	ASN	C-N-CA	-5.60	110.54	122.30
1	C	839	ASP	C-N-CA	-5.59	107.72	121.70
1	A	1098	ASN	C-N-CA	-5.46	110.83	122.30
1	C	557	LYS	C-N-CA	5.38	135.16	121.70
1	A	351	TYR	C-N-CA	5.32	135.01	121.70
1	C	841	LEU	CA-C-N	5.31	126.82	116.20
1	B	406	GLU	C-N-CA	5.27	134.87	121.70
1	B	462	LYS	CA-C-N	-5.25	102.39	117.10
1	B	515	PHE	N-CA-C	-5.21	96.95	111.00
1	B	515	PHE	O-C-N	5.19	131.01	122.70
1	C	351	TYR	C-N-CA	5.19	134.67	121.70
1	C	700	GLY	N-CA-C	-5.17	100.18	113.10
1	A	174	PRO	N-CA-C	5.15	125.49	112.10
1	B	854	LYS	CA-C-N	-5.06	106.06	117.20
1	B	585	LEU	CA-C-O	-5.05	109.49	120.10
1	B	351	TYR	C-N-CA	5.02	134.26	121.70
1	C	841	LEU	C-N-CA	5.01	132.81	122.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	586	ASP	Peptide
1	B	854	LYS	Peptide

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4484	0	1220	7	0
1	B	4484	0	1220	24	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	4484	0	1219	22	0
All	All	13452	0	3659	34	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 (34) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:586:ASP:N	1:C:842:GLY:HA3	1.90	0.85
1:B:586:ASP:CA	1:C:844:ILE:H	1.93	0.81
1:B:199:GLY:HA2	1:B:232:GLY:HA2	1.61	0.81
1:B:199:GLY:CA	1:B:232:GLY:HA2	2.19	0.73
1:B:586:ASP:CA	1:C:843:ASP:N	2.53	0.72
1:A:136:CYS:H	1:A:159:VAL:CA	2.04	0.71
1:B:585:LEU:C	1:C:842:GLY:HA3	2.12	0.68
1:B:575:ALA:CA	1:C:843:ASP:CA	2.72	0.68
1:B:199:GLY:C	1:B:232:GLY:HA2	2.15	0.68
1:C:729:VAL:H	1:C:1059:GLY:HA2	1.62	0.64
1:B:587:ILE:CA	1:C:842:GLY:N	2.64	0.60
1:B:587:ILE:N	1:C:842:GLY:C	2.57	0.57
1:B:586:ASP:C	1:C:844:ILE:H	2.09	0.56
1:C:142:GLY:HA3	1:C:244:LEU:O	2.07	0.55
1:B:585:LEU:O	1:C:843:ASP:N	2.41	0.54
1:B:587:ILE:N	1:C:843:ASP:N	2.55	0.54
1:B:586:ASP:CA	1:C:843:ASP:CA	2.90	0.50
1:A:44:ARG:O	1:A:283:GLY:HA2	2.13	0.49
1:B:729:VAL:H	1:B:1059:GLY:HA2	1.79	0.47
1:A:138:ASP:C	1:A:140:PHE:H	2.18	0.46
1:B:586:ASP:O	1:C:841:LEU:C	2.54	0.45
1:B:148:ASN:C	1:B:150:LYS:H	2.20	0.45
1:A:320:VAL:CA	1:A:623:ALA:CA	2.94	0.45
1:B:553:THR:N	1:C:841:LEU:O	2.50	0.44
1:A:109:THR:C	1:A:111:ASP:H	2.22	0.42
1:B:586:ASP:CA	1:C:844:ILE:N	2.74	0.42
1:A:729:VAL:H	1:A:1059:GLY:HA2	1.83	0.42
1:C:698:SER:C	1:C:700:GLY:H	2.21	0.42
1:A:145:TYR:N	1:A:246:ARG:O	2.52	0.42
1:B:588:THR:N	1:C:841:LEU:N	2.68	0.42
1:B:551:VAL:O	1:C:841:LEU:CA	2.68	0.41
1:B:553:THR:H	1:C:841:LEU:C	2.24	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:586:ASP:C	1:C:843:ASP:N	2.74	0.41
1:B:586:ASP:O	1:C:840:CYS:O	2.39	0.41

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	A	1119/1249 (90%)	1054 (94%)	49 (4%)	16 (1%)	11	44
1	B	1119/1249 (90%)	1032 (92%)	64 (6%)	23 (2%)	7	36
1	C	1119/1249 (90%)	1046 (94%)	58 (5%)	15 (1%)	12	46
All	All	3357/3747 (90%)	3132 (93%)	171 (5%)	54 (2%)	13	42

All (54) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	THR
1	A	353	TRP
1	A	689	SER
1	A	796	ASP
1	B	40	ASP
1	B	255	SER
1	B	412	PRO
1	B	467	ASP
1	B	516	GLU
1	B	689	SER
1	B	796	ASP
1	C	153	MET
1	C	796	ASP
1	C	841	LEU
1	A	158	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	255	SER
1	A	794	ILE
1	B	375	SER
1	B	429	PHE
1	B	687	VAL
1	B	744	GLY
1	C	75	GLY
1	C	353	TRP
1	C	707	TYR
1	C	833	PHE
1	A	76	THR
1	A	378	LYS
1	A	429	PHE
1	A	690	GLN
1	A	707	TYR
1	A	797	PHE
1	B	353	TRP
1	B	571	ASP
1	B	813	SER
1	C	429	PHE
1	C	1098	ASN
1	A	1098	ASN
1	B	155	SER
1	B	414	GLN
1	B	707	TYR
1	B	721	SER
1	C	146	HIS
1	C	534	VAL
1	C	688	ALA
1	B	402	ILE
1	B	410	ILE
1	B	440	ASN
1	B	987	PRO
1	C	813	SER
1	B	411	ALA
1	A	1065	VAL
1	A	381	GLY
1	C	476	GLY
1	C	173	GLN

### 5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

### 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](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 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.

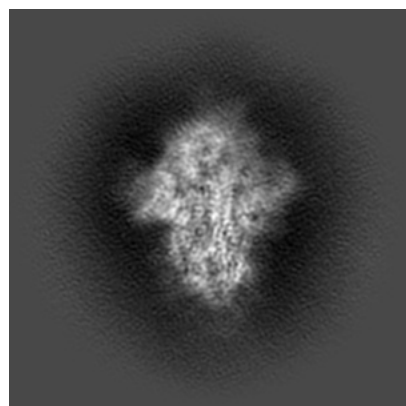
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-30915. These allow visual inspection of the internal detail of the map and identification of artifacts.

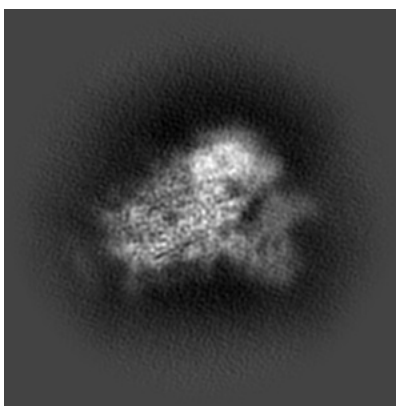
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

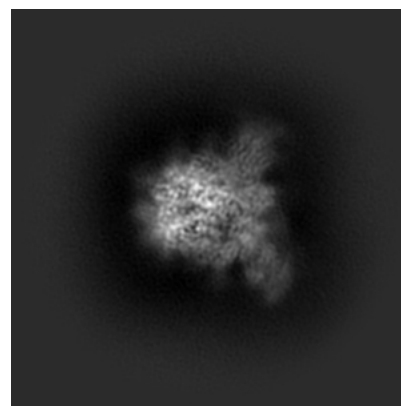
#### 6.1.1 Primary map



X

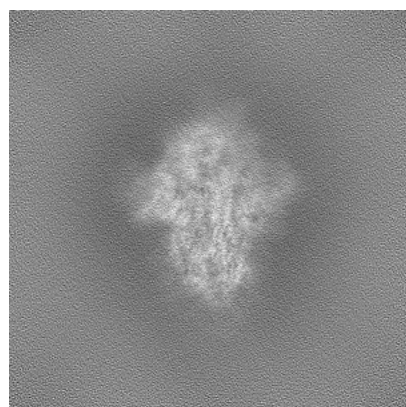


Y

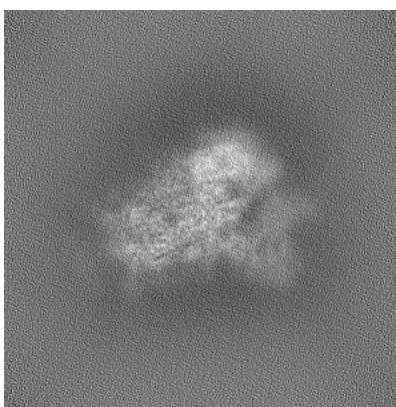


Z

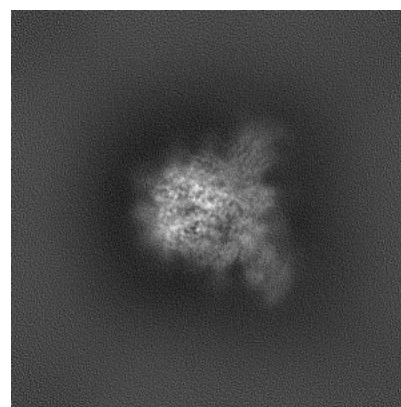
#### 6.1.2 Raw map



X



Y



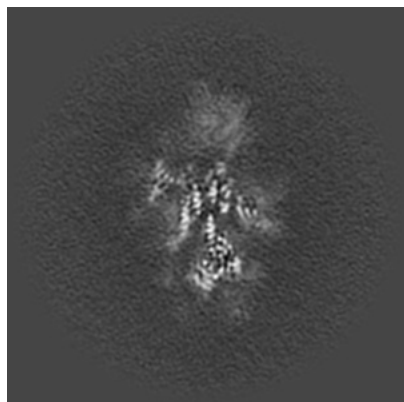
Z

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

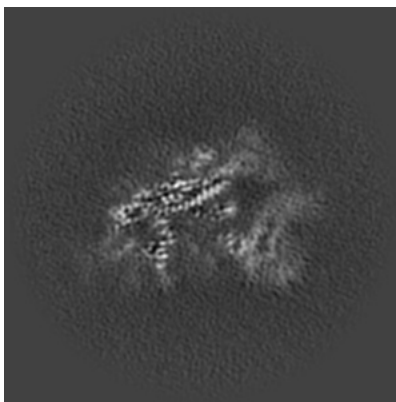


## 6.2 Central slices [i](#)

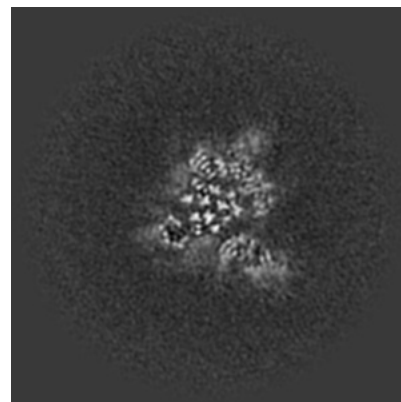
### 6.2.1 Primary map



X Index: 180

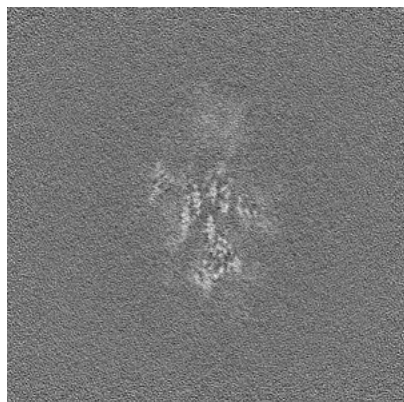


Y Index: 180

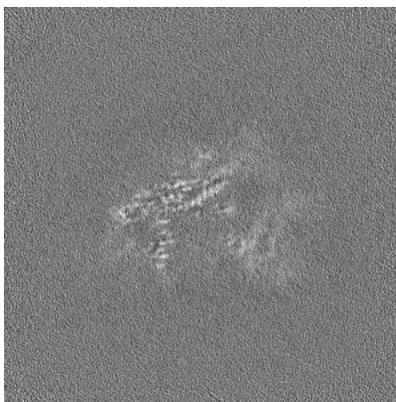


Z Index: 180

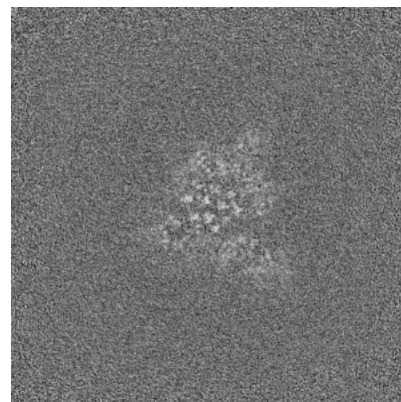
### 6.2.2 Raw map



X Index: 180



Y Index: 180



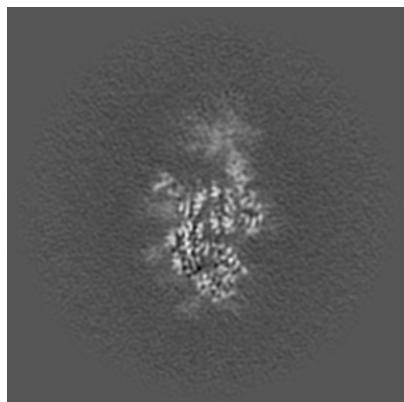
Z Index: 180

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

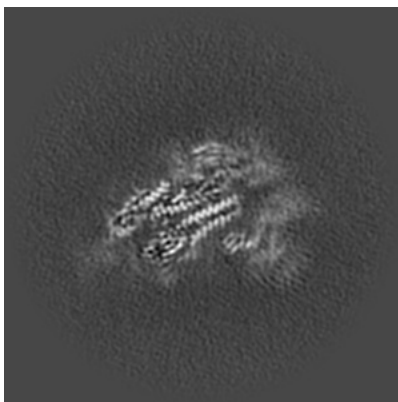


## 6.3 Largest variance slices [i](#)

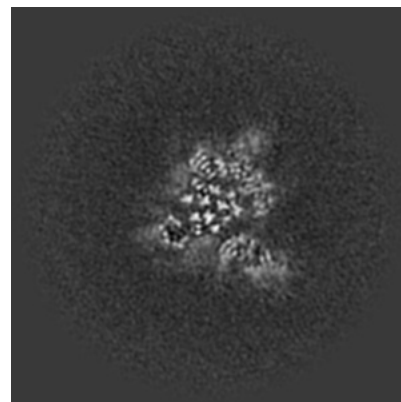
### 6.3.1 Primary map



X Index: 167

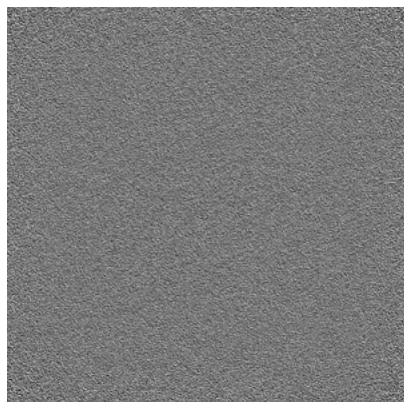


Y Index: 185

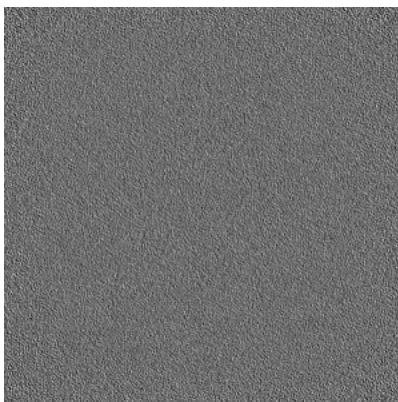


Z Index: 180

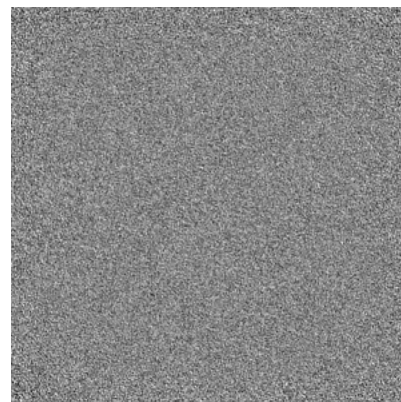
### 6.3.2 Raw map



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 surface views [i](#)

### 6.4.1 Primary map



X



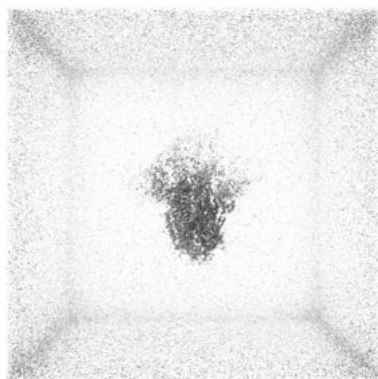
Y



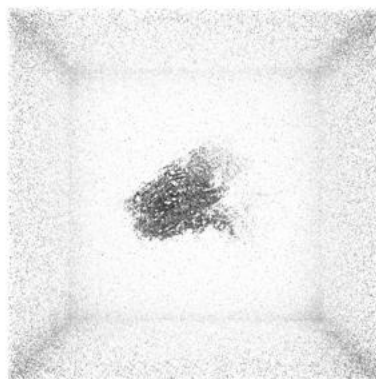
Z

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.

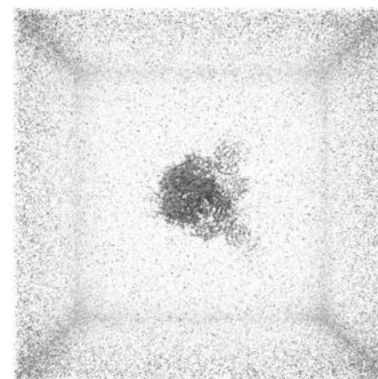
### 6.4.2 Raw map



X



Y



Z

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.

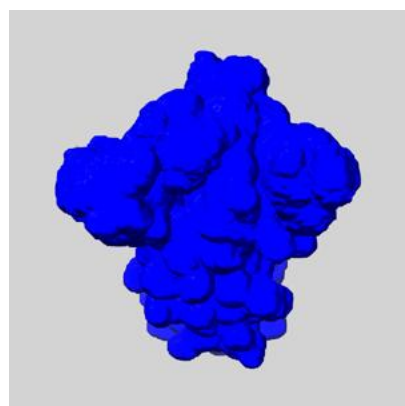
## 6.5 Mask visualisation [i](#)

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

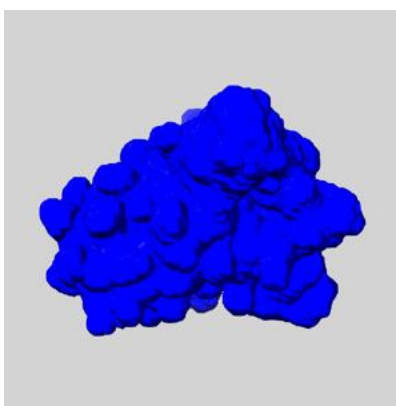
A mask typically either:

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

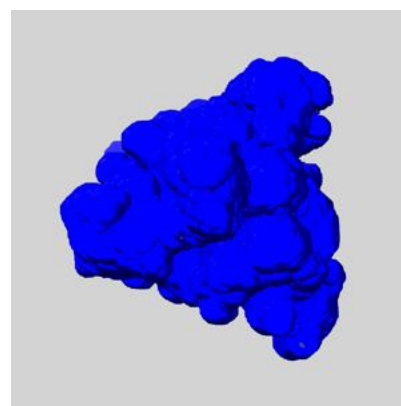
### 6.5.1 emd\_30915\_msk\_1.map [i](#)



X



Y

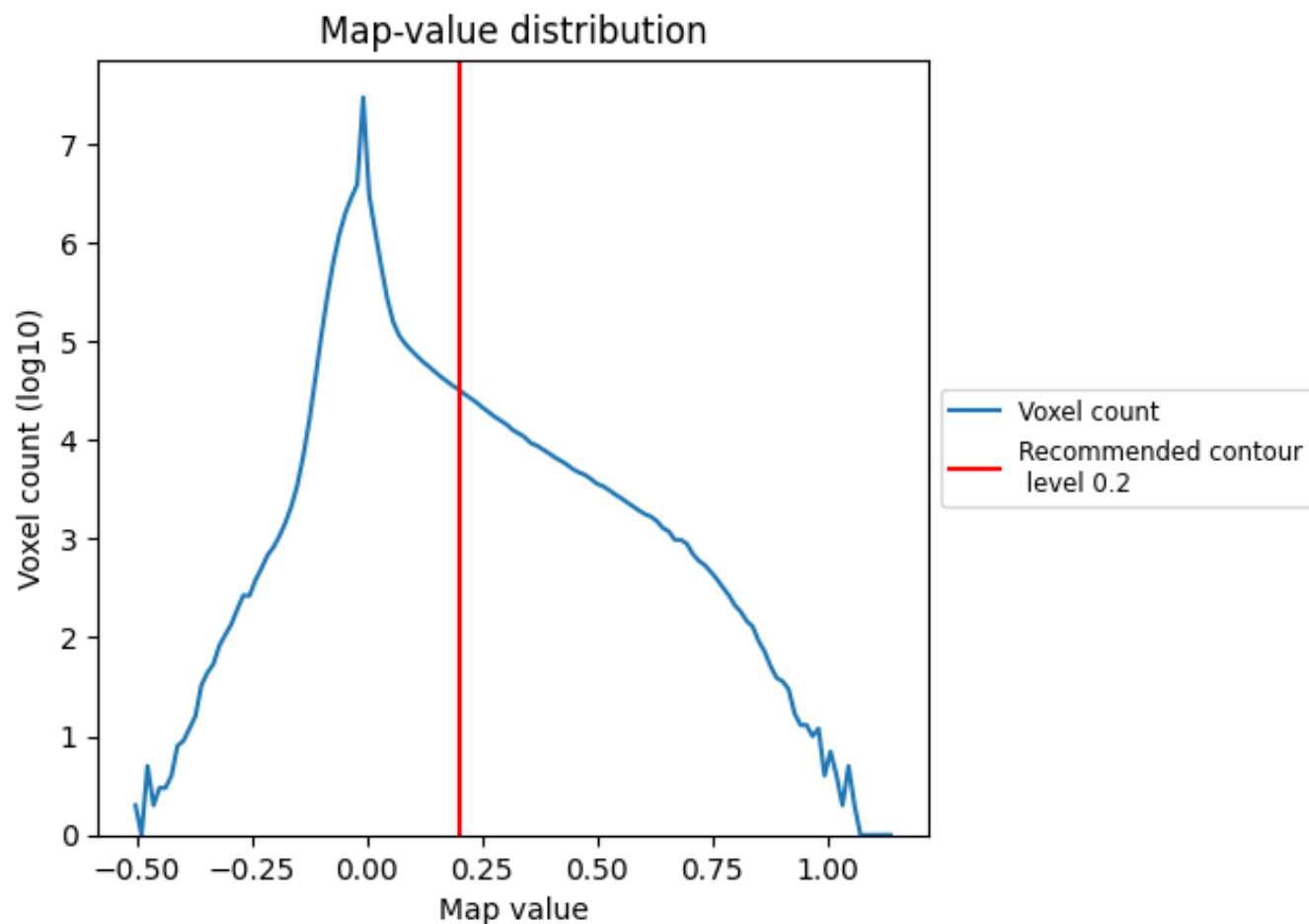


Z

## 7 Map analysis [i](#)

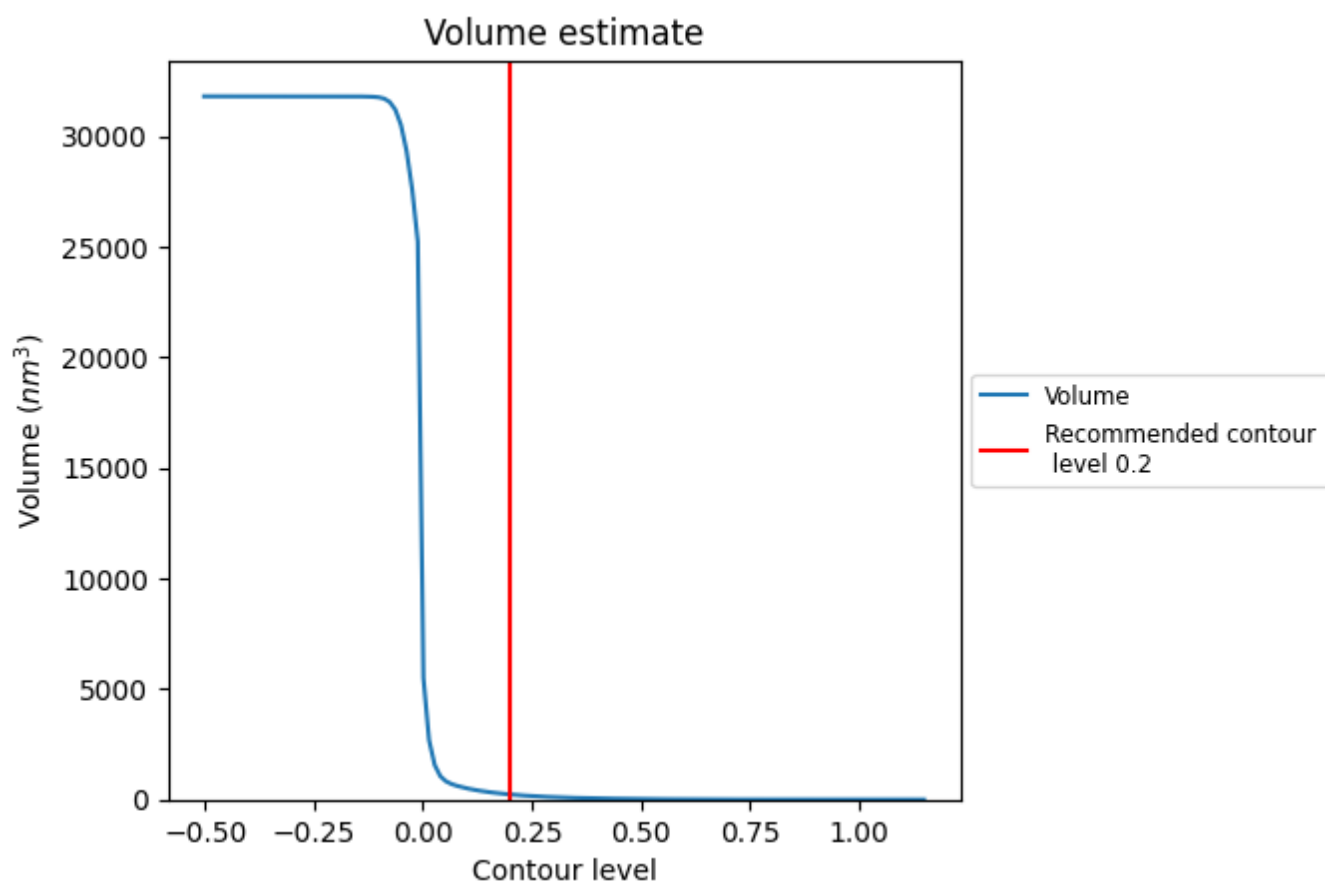
This section contains the results of statistical analysis of the map.

### 7.1 Map-value distribution [i](#)



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

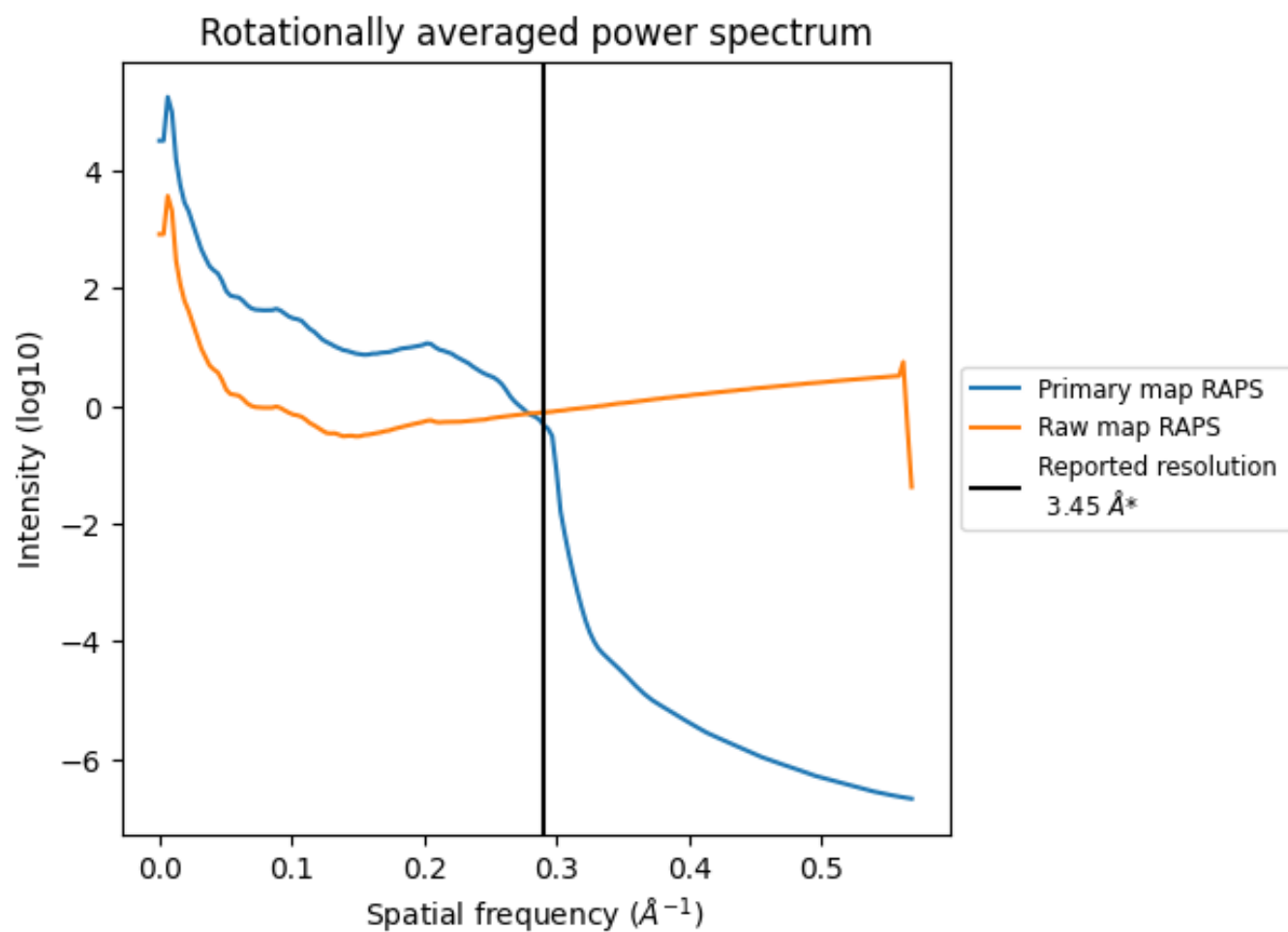
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 232 nm<sup>3</sup>; this corresponds to an approximate mass of 209 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 ⓘ

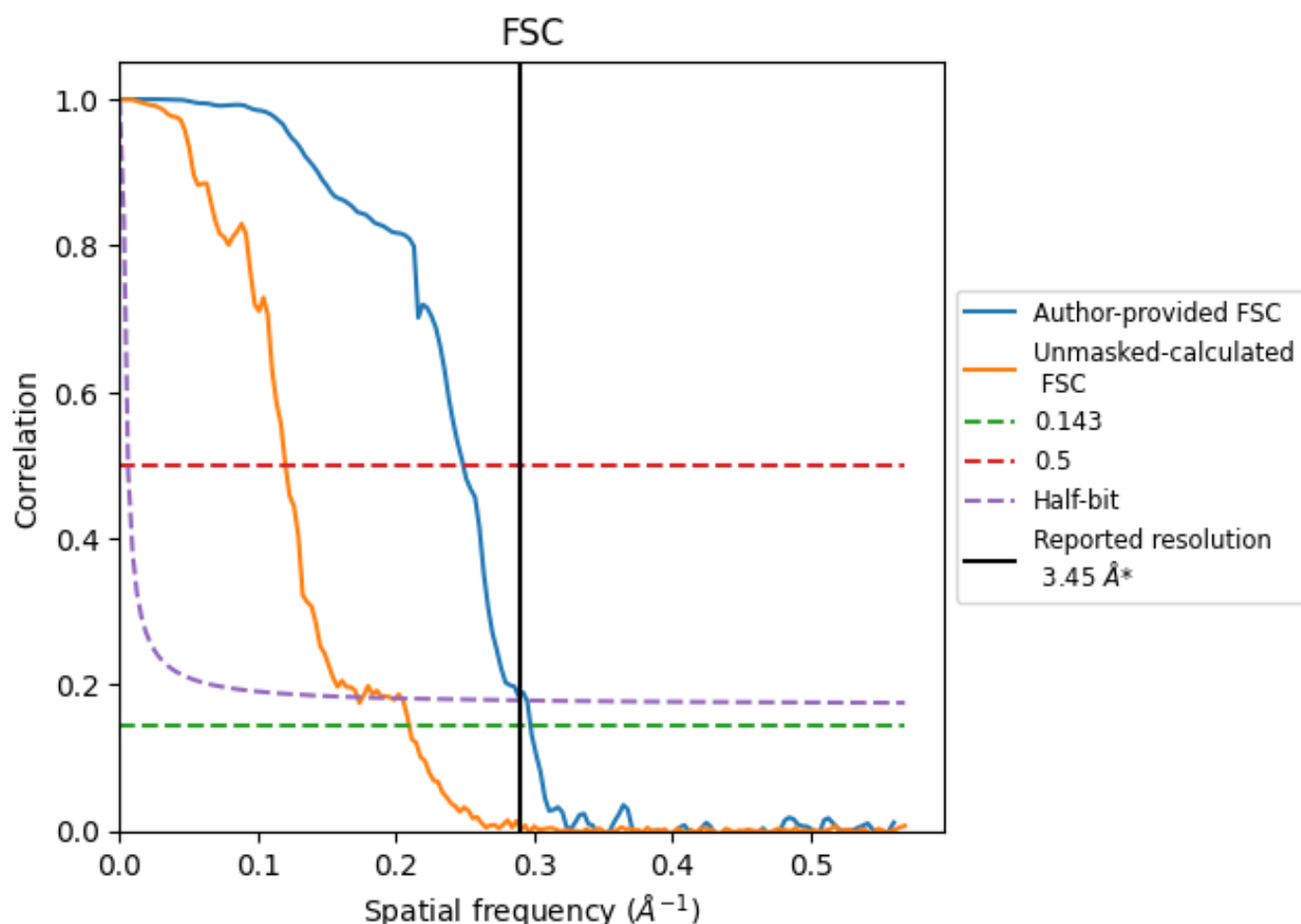


\*Reported resolution corresponds to spatial frequency of 0.290  $\text{\AA}^{-1}$

## 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 0.290 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.45	-	-
Author-provided FSC curve	3.36	4.02	3.39
Unmasked-calculated*	4.77	8.33	5.81

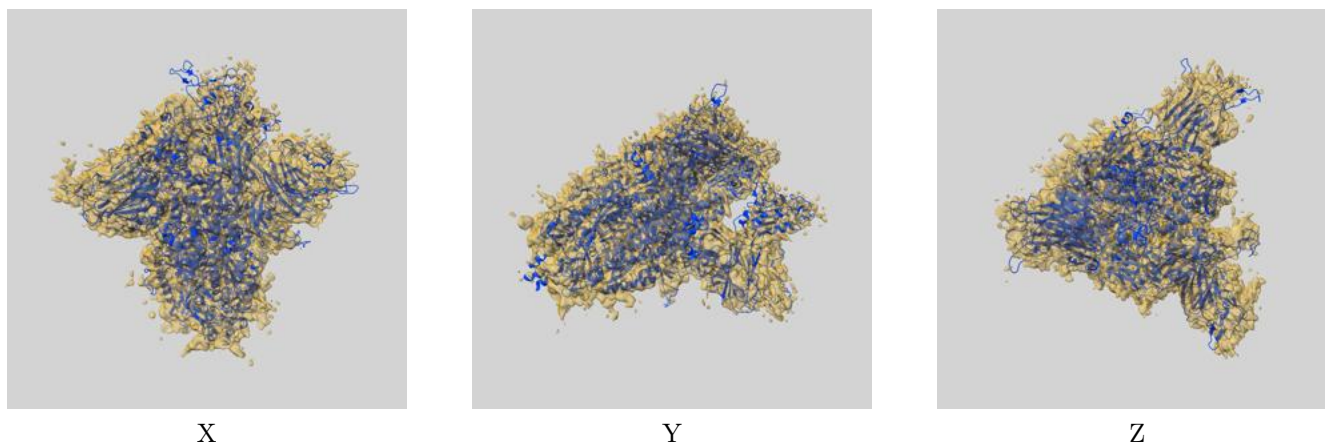
\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.77 differs from the reported value 3.45 by more than 10 %



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

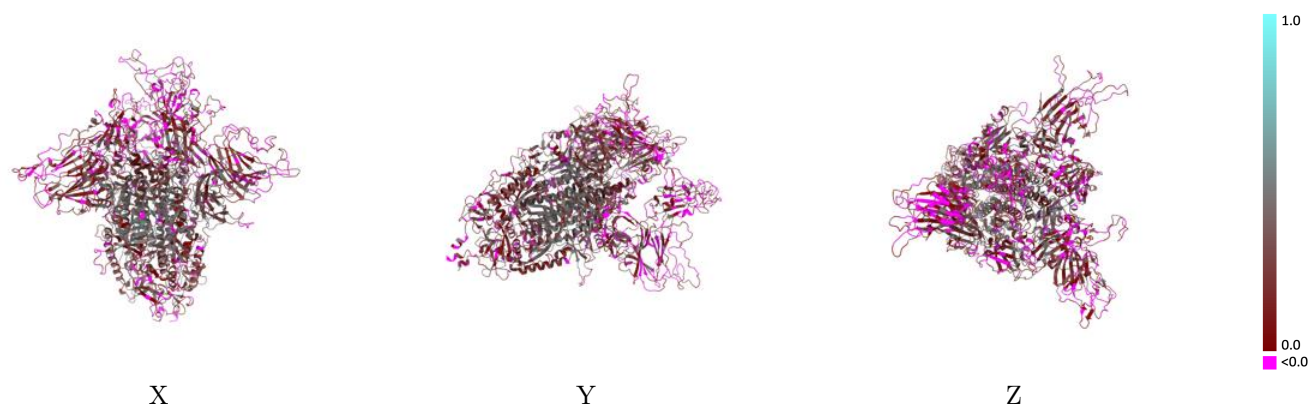
This section contains information regarding the fit between EMDB map EMD-30915 and PDB model 7DZW. Per-residue inclusion information can be found in [section 3](#) on [page 5](#).

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



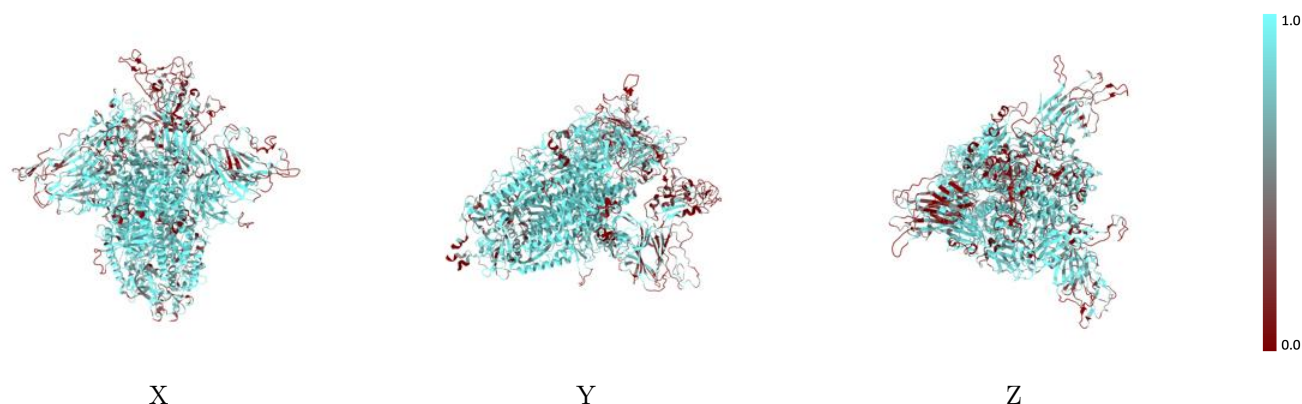
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



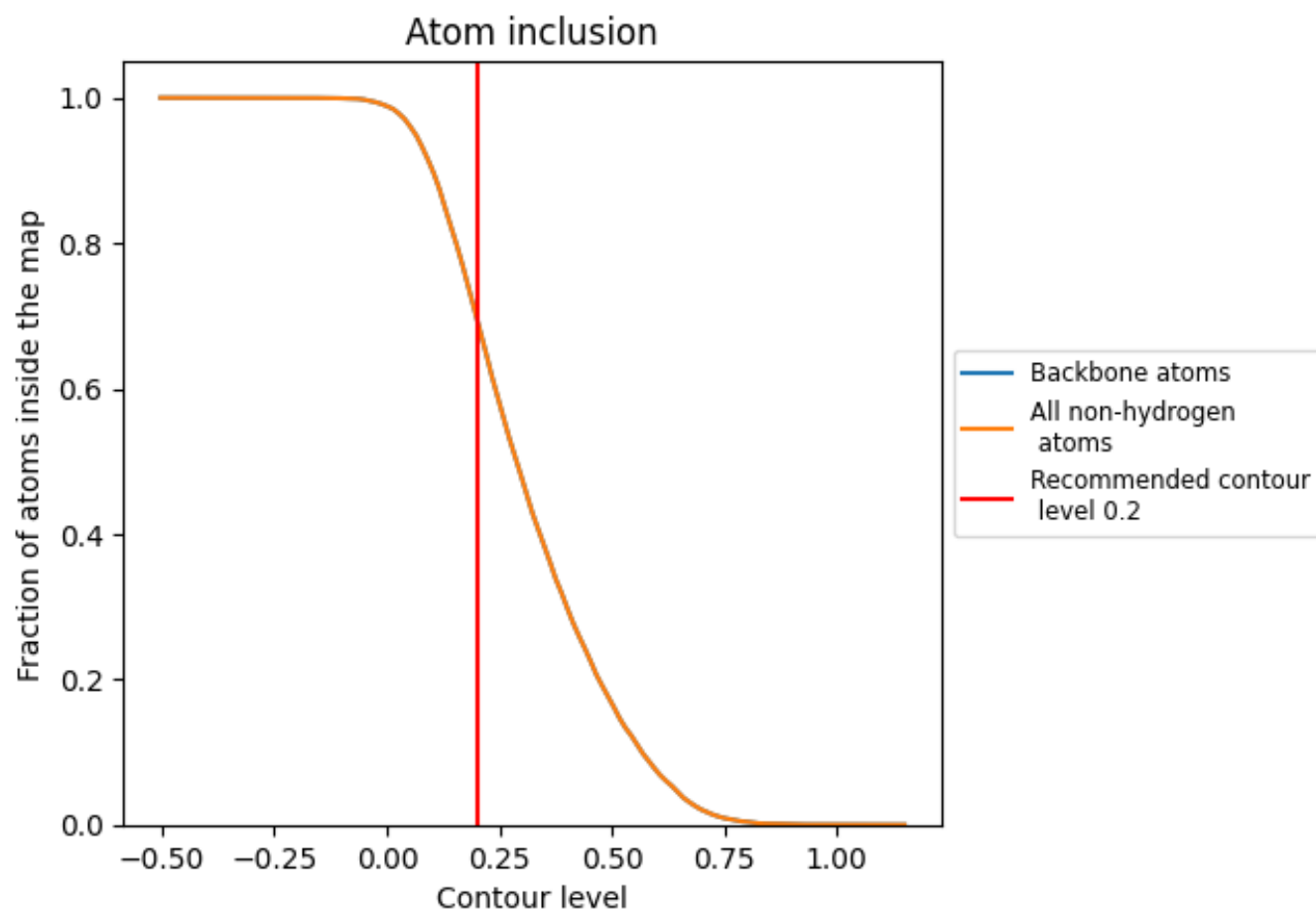
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 70% of all backbone atoms, 70% 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	<div></div> 0.6965	<div></div> 0.2060
A	<div></div> 0.7244	<div></div> 0.2240
B	<div></div> 0.6579	<div></div> 0.2040
C	<div></div> 0.7072	<div></div> 0.1910

