

wwPDB EM Validation Summary Report (i)

Mar 12, 2024 – 02:52 PM JST

PDB ID : 8JHZ

EMDB ID : EMD-36301

Title : Cryo-EM structure of the TcsH-TMPRSS2 complex

Authors: Zhou, R.; Liang, T.; Zhan, X.

Deposited on : 2023-05-25

Resolution : 3.20 Å(reported)

This is a wwPDB EM Validation Summary 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 (i) 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 (i)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev70

Mogul : 1.8.5 (274361), CSD as541be (2020)

MolProbity: 4.02b-467

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $MapQ \quad : \quad 1.9.13$

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

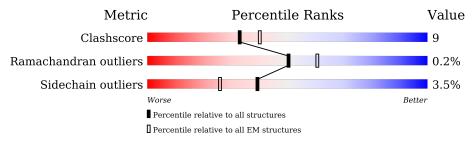
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $ELECTRON\ MICROSCOPY$

The reported resolution of this entry is 3.20 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 $(\# \mathrm{Entries})$	${ m EM\ structures} \ (\#{ m Entries})$
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 <=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	2626	7% 81%			17%	.
2	В	424	55% 22°	2%		20%	_



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 23814 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 Hemorrhagic toxin.

\mathbf{Mol}	Chain	Residues	Atoms						AltConf	Trace
1	A	2615	Total 21144	C 13557	N 3425	O 4120	P 1	S 41	0	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1833	SER	LEU	conflict	UNP M9ZTT7
A	2619	HIS	-	expression tag	UNP M9ZTT7
A	2620	HIS	-	expression tag	UNP M9ZTT7
A	2621	HIS	-	expression tag	UNP M9ZTT7
A	2622	HIS	-	expression tag	UNP M9ZTT7
A	2623	HIS	-	expression tag	UNP M9ZTT7
A	2624	HIS	-	expression tag	UNP M9ZTT7
A	2625	HIS	-	expression tag	UNP M9ZTT7
A	2626	HIS	-	expression tag	UNP M9ZTT7

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

Mol	Chain	Residues	Atoms				AltConf	Trace	
2	В	341	Total 2669	C 1688	N 462	O 496	S 23	0	0

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	69	MET	-	initiating methionine	UNP O15393
В	70	GLY	-	expression tag	UNP O15393
В	71	ILE	-	expression tag	UNP O15393
В	72	LEU	-	expression tag	UNP O15393
В	73	PRO	-	expression tag	UNP O15393
В	74	SER	-	expression tag	UNP O15393
В	75	PRO	-	expression tag	UNP O15393
В	76	GLY	-	expression tag	UNP O15393

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Chain	Residue	Modelled	Actual	Comment	Reference
В	77	MET	-	expression tag	UNP O15393
В	78	PRO	-	expression tag	UNP O15393
В	79	ALA	-	expression tag	UNP O15393
В	80	LEU	-	expression tag	UNP O15393
В	81	LEU	-	expression tag	UNP O15393
В	82	SER	-	expression tag	UNP O15393
В	83	LEU	-	expression tag	UNP O15393
В	84	VAL	-	expression tag	UNP O15393
В	85	SER	-	expression tag	UNP O15393
В	86	LEU	-	expression tag	UNP O15393
В	87	LEU	-	expression tag	UNP O15393
В	88	SER	-	expression tag	UNP O15393
В	89	VAL	-	expression tag	UNP O15393
В	90	LEU	-	expression tag	UNP O15393
В	91	LEU	-	expression tag	UNP O15393
В	92	MET	-	expression tag	UNP O15393
В	93	GLY	-	expression tag	UNP O15393
В	94	CYS	-	expression tag	UNP O15393
В	95	VAL	-	expression tag	UNP O15393
В	96	ALA	-	expression tag	UNP O15393
В	97	GLU	-	expression tag	UNP O15393
В	98	THR	-	expression tag	UNP O15393
В	99	GLY	-	expression tag	UNP O15393
В	100	HIS	-	expression tag	UNP O15393
В	101	HIS	-	expression tag	UNP O15393
В	102	HIS	-	expression tag	UNP O15393
В	103	HIS	-	expression tag	UNP O15393
В	104	HIS	-	expression tag	UNP O15393
В	105	HIS	-	expression tag	UNP O15393
В	255	GLN	ARG	conflict	UNP O15393

 \bullet Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

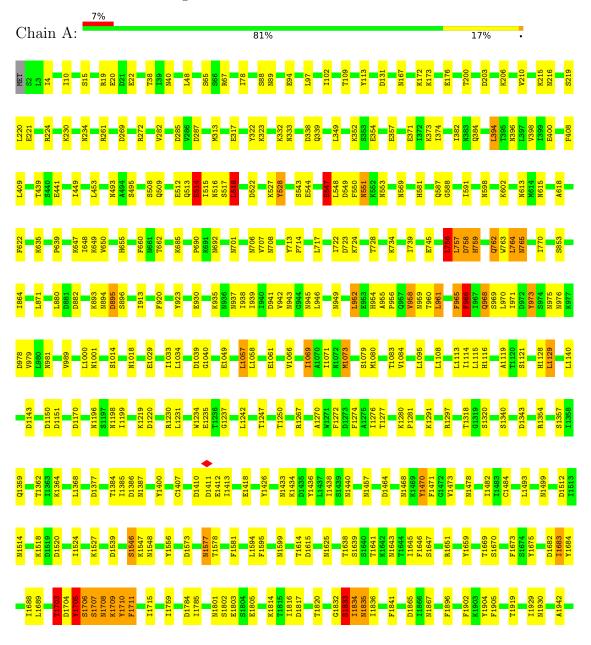
Mol	Chain	Residues	Atoms	AltConf
3	A	1	Total Zn 1 1	0



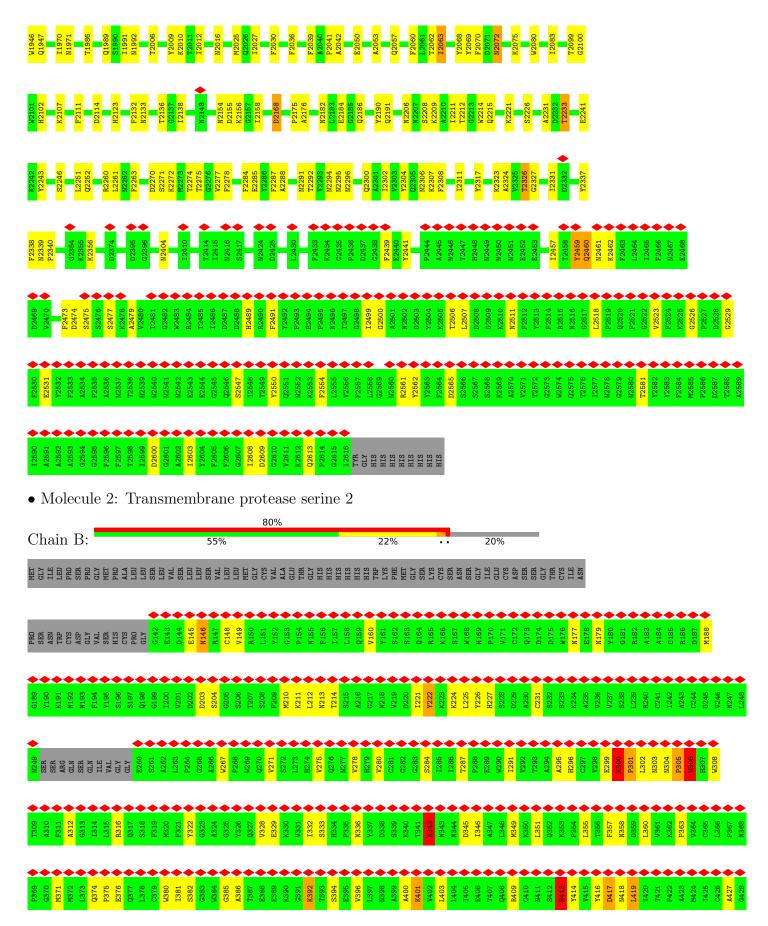
3 Residue-property plots (i)

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

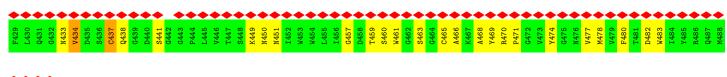
• Molecule 1: Hemorrhagic toxin















4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	87378	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 $(e^-/\text{Å}^2)$	50.00	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	5.558	Depositor
Minimum map value	-2.531	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.065	Depositor
Recommended contour level	0.35	Depositor
Map size (Å)	652.2, 652.2, 652.2	wwPDB
Map dimensions	600, 600, 600	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.087, 1.087, 1.087	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: ZN, SEP

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	Bo	ond lengths	Bond angles		
MIOI	RMSZ # Z >		# Z > 5	RMSZ	# Z >5	
1	A	0.50	$22/21600 \ (0.1\%)$	0.57	$22/29255 \ (0.1\%)$	
2	В	0.55	8/2740 (0.3%)	0.74	13/3724 (0.3%)	
All	All	0.51	30/24340 (0.1%)	0.59	$35/32979 \ (0.1\%)$	

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 maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
2	В	0	2
All	All	0	5

The worst 5 of 30 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	Ideal(Å)
2	В	304	ASN	C-O	-9.72	1.04	1.23
1	A	1833	SER	C-O	-9.71	1.04	1.23
1	A	1832	GLY	C-O	-9.40	1.08	1.23
1	A	1471	PHE	C-O	-9.06	1.06	1.23
1	A	1710	TYR	C-O	-8.11	1.07	1.23

The worst 5 of 35 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
1	A	408	PHE	CB-CA-C	-14.12	82.17	110.40
1	A	959	ASN	CB-CA-C	-10.30	89.80	110.40
1	A	1471	PHE	CB-CA-C	-9.20	92.01	110.40
1	A	960	THR	CA-CB-OG1	-8.96	90.19	109.00

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Mol	Chain	Res	Type	Atoms	${f Z}$	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	A	1835	ASN	CB-CA-C	-8.86	92.69	110.40

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1703	THR	Mainchain
1	A	1833	SER	Mainchain
1	A	756	LEU	Mainchain
2	В	300	LYS	Mainchain
2	В	306	TRP	Mainchain

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	A	21144	0	20628	374	0
2	В	2669	0	2561	58	0
3	A	1	0	0	0	0
All	All	23814	0	23189	431	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 9.

The worst 5 of 431 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:2308:PHE:CZ	1:A:2331:ILE:HD13	1.37	1.56
1:A:2326:THR:CB	1:A:2340:PRO:HA	1.49	1.42
1:A:2308:PHE:HZ	1:A:2331:ILE:CD1	1.32	1.39
1:A:2326:THR:OG1	1:A:2340:PRO:CA	1.75	1.33
1:A:655:HIS:HE1	1:A:759:HIS:CE1	1.51	1.28

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	Analysed Favoured Allowed		Outliers	Perce	\mathbf{ntiles}
1	A	$2612/2626\ (100\%)$	2499 (96%)	109 (4%)	4 (0%)	47	79
2	В	337/424~(80%)	300 (89%)	34 (10%)	3 (1%)	17	56
All	All	$2949/3050 \ (97\%)$	2799 (95%)	143 (5%)	7 (0%)	50	79

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1707	SER
2	В	300	LYS
1	A	1703	THR
2	В	145	GLU
1	A	1705	TYR

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	A	2352/2362 (100%)	2290 (97%)	62 (3%)	46 76	
2	В	290/360 (81%)	259 (89%)	31 (11%)	6 27	
All	All	2642/2722 (97%)	2549 (96%)	93 (4%)	39 69	

5 of 93 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2460	GLN

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Mol	Chain	Res	Type
2	В	306	TRP
1	A	2565	ASP
2	В	204	SER
2	В	342	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 41 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1909	ASN
1	A	2047	ASN
1	A	1959	ASN
1	A	1993	ASN
1	A	2058	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)

1 non-standard protein/DNA/RNA residue 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		ype Chain Res Link		B	Bond lengths		Bond angles			
Moi Type	Chain	rtes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
1	SEP	A	517	1	8,9,10	1.54	2 (25%)	8,12,14	2.01	2 (25%)

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
1	SEP	A	517	1	-	4/5/8/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$\operatorname{Ideal}(ext{\AA})$
1	A	517	SEP	P-OG	2.40	1.67	1.60
1	A	517	SEP	P-O3P	-2.22	1.46	1.54

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
1	A	517	SEP	OG-P-O1P	3.87	117.34	106.47
1	A	517	SEP	OG-CB-CA	2.95	111.02	108.14

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	517	SEP	N-CA-CB-OG
1	A	517	SEP	CB-OG-P-O2P
1	A	517	SEP	CB-OG-P-O3P
1	A	517	SEP	CB-OG-P-O1P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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.



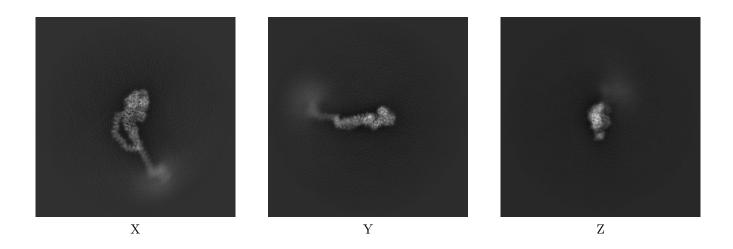
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-36301. These allow visual inspection of the internal detail of the map and identification of artifacts.

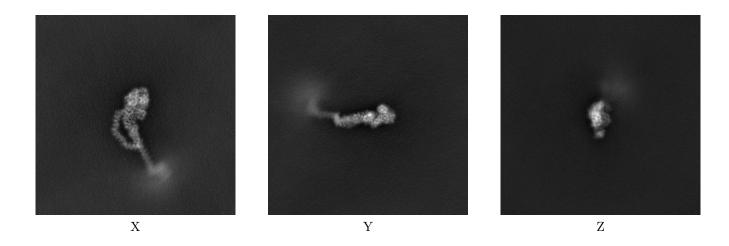
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

6.1.1 Primary map



6.1.2 Raw map



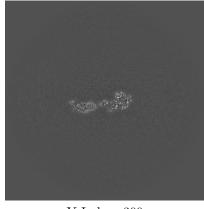
The images above show the map projected in three orthogonal directions.



6.2 Central slices (i)

6.2.1 Primary map







X Index: 300

Y Index: 300

Z Index: 300

6.2.2 Raw map







X Index: 300

Y Index: 300

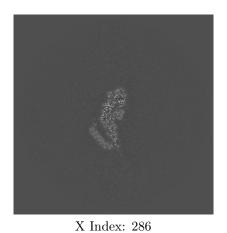
Z Index: 300

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



6.3 Largest variance slices (i)

Primary map 6.3.1





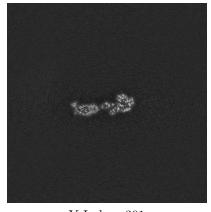


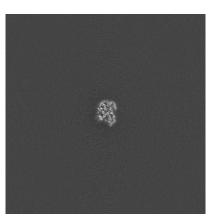
Y Index: 292

Z Index: 340

Raw map 6.3.2







X Index: 285

Y Index: 301

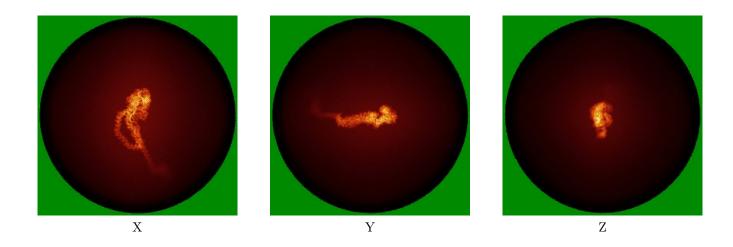
Z Index: 340

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

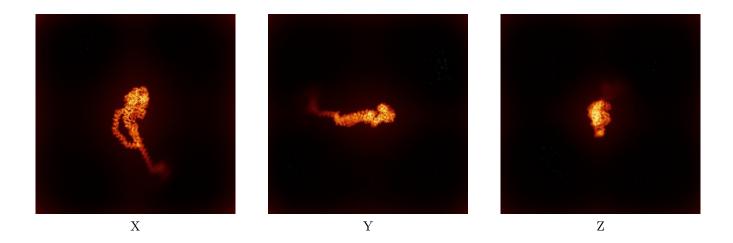


6.4 Orthogonal standard-deviation projections (False-color) (i)

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 0.35. 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)

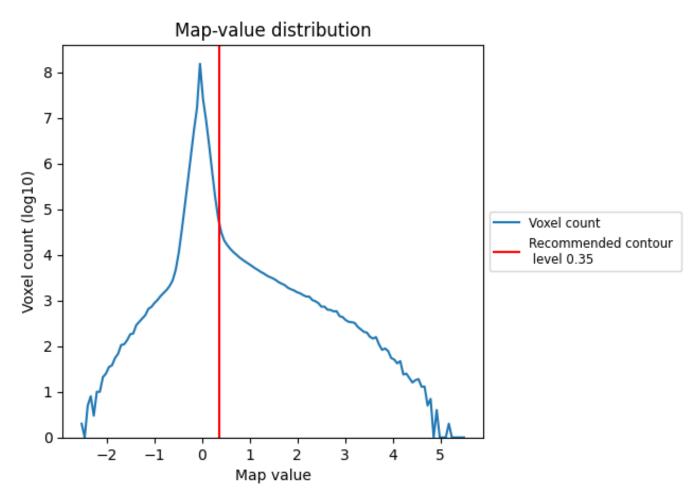
This section was not generated. No masks/segmentation were deposited.



7 Map analysis (i)

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

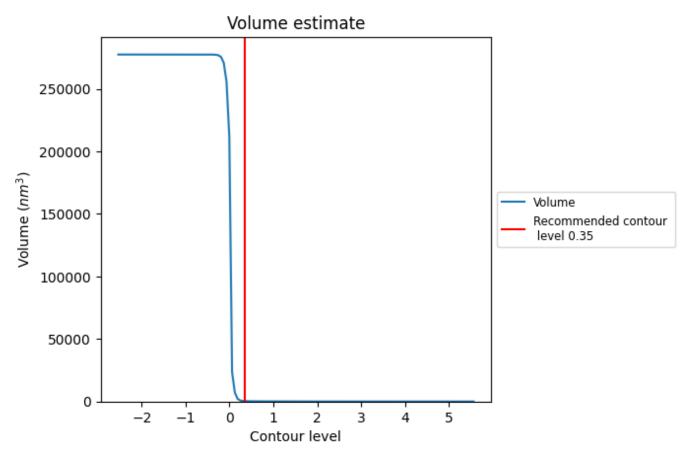
7.1 Map-value distribution (i)



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)

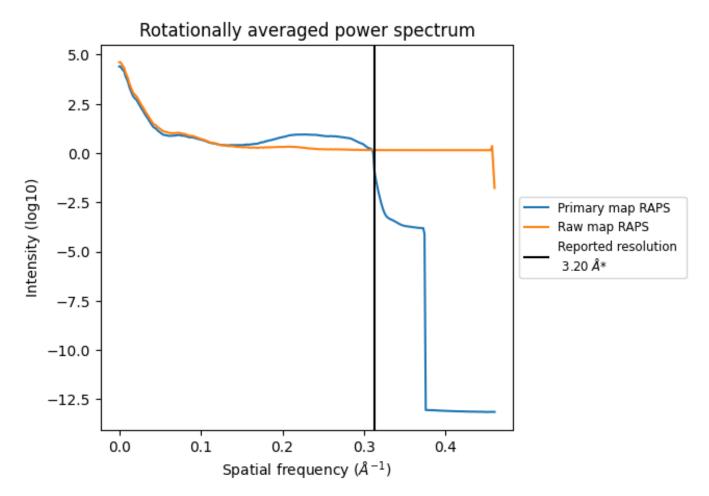


The volume at the recommended contour level is $325~\mathrm{nm}^3$; this corresponds to an approximate mass of $293~\mathrm{kDa}$.

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)



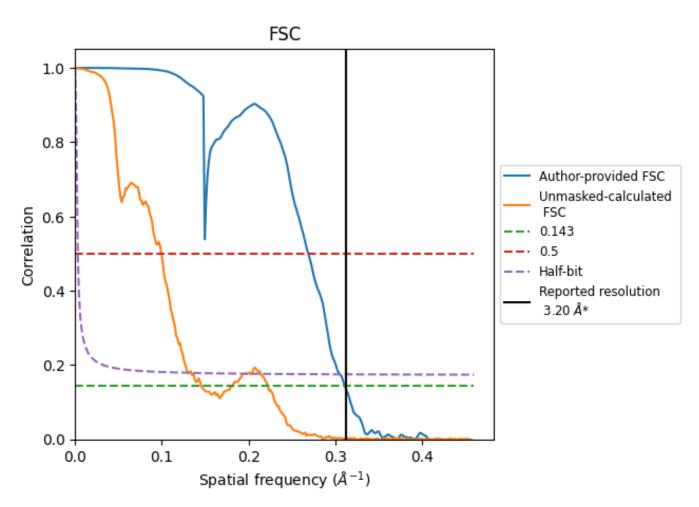
*Reported resolution corresponds to spatial frequency of 0.312 $\rm \mathring{A}^{-1}$



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 0.312 $\rm \mathring{A}^{-1}$



8.2 Resolution estimates (i)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)			
rtesolution estimate (A)	0.143	0.5	Half-bit	
Reported by author	3.20	-	-	
Author-provided FSC curve	3.21	3.72	3.28	
Unmasked-calculated*	6.87	10.06	7.67	

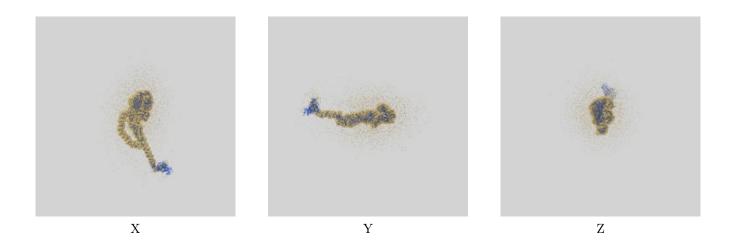
^{*}Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 6.87 differs from the reported value 3.2 by more than 10 %



9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-36301 and PDB model 8JHZ. Per-residue inclusion information can be found in section 3 on page 5.

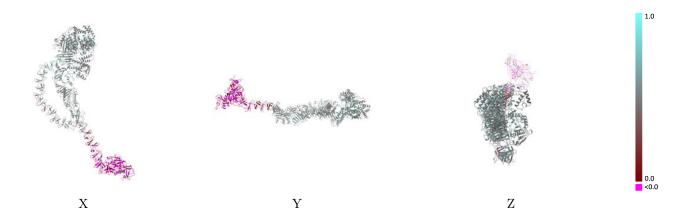
9.1 Map-model overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.35 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.

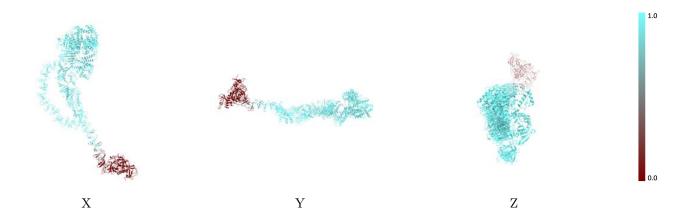


9.2 Q-score mapped to coordinate model (i)



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.

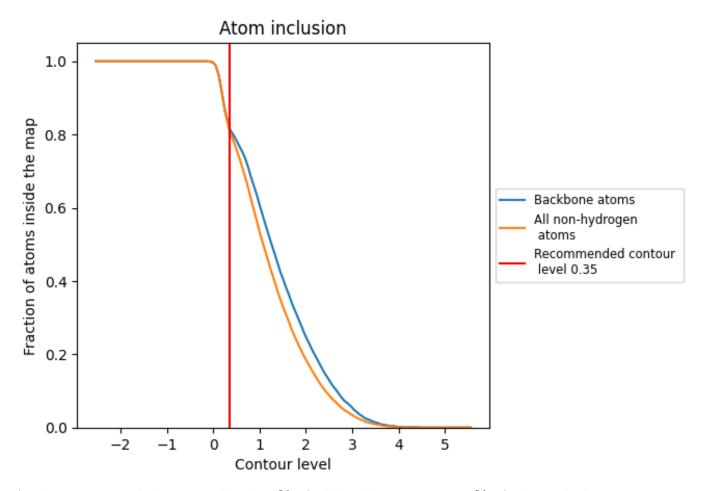
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.35).



9.4 Atom inclusion (i)



At the recommended contour level, 82% of all backbone atoms, 81% of all non-hydrogen atoms, are inside the map.



9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.35) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8110	0.4040
A	0.9100	0.4560
В	0.0210	-0.0020



