



wwPDB EM Validation Summary Report ⓘ

Feb 10, 2022 – 01:08 pm GMT

PDB ID : 7PG4
EMDB ID : EMD-13388
Title : Low resolution Cryo-EM structure of the full-length insulin receptor bound to 2 insulin, conf 3
Authors : Nielsen, J.A.; Slaaby, R.; Boesen, T.; Hummelshoj, T.; Brandt, J.; Schluckebier, G.; Nissen, P.
Deposited on : 2021-08-12
Resolution : 9.10 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 ⓘ symbol.

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

EMDB validation analysis : 0.0.0.dev97
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.26

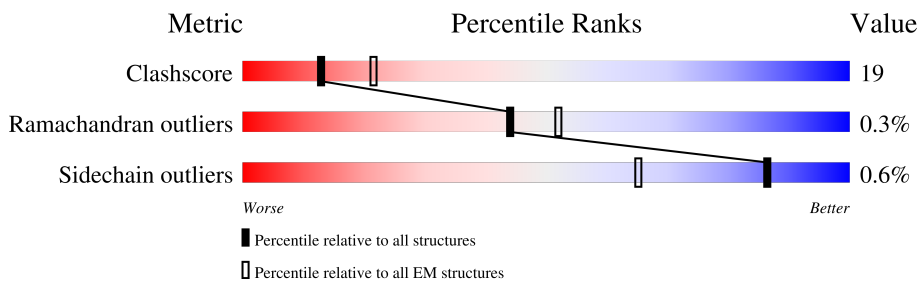
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 9.10 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	A	1382	
1	B	1382	
2	C	21	
2	I	21	
3	D	30	
3	J	30	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 13682 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 Isoform Short of Insulin receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	786	6340	4014	1099	1178	49	0	0
1	B	822	6631	4201	1139	1239	52	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1344	GLU	-	expression tag	UNP P06213
A	1345	ASP	-	expression tag	UNP P06213
A	1346	GLN	-	expression tag	UNP P06213
A	1347	VAL	-	expression tag	UNP P06213
A	1348	ASP	-	expression tag	UNP P06213
A	1349	PRO	-	expression tag	UNP P06213
A	1350	ARG	-	expression tag	UNP P06213
A	1351	LEU	-	expression tag	UNP P06213
A	1352	ILE	-	expression tag	UNP P06213
A	1353	ASP	-	expression tag	UNP P06213
A	1354	GLY	-	expression tag	UNP P06213
A	1355	LYS	-	expression tag	UNP P06213
B	1344	GLU	-	expression tag	UNP P06213
B	1345	ASP	-	expression tag	UNP P06213
B	1346	GLN	-	expression tag	UNP P06213
B	1347	VAL	-	expression tag	UNP P06213
B	1348	ASP	-	expression tag	UNP P06213
B	1349	PRO	-	expression tag	UNP P06213
B	1350	ARG	-	expression tag	UNP P06213
B	1351	LEU	-	expression tag	UNP P06213
B	1352	ILE	-	expression tag	UNP P06213
B	1353	ASP	-	expression tag	UNP P06213
B	1354	GLY	-	expression tag	UNP P06213
B	1355	LYS	-	expression tag	UNP P06213

- Molecule 2 is a protein called Insulin.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	21	Total	C	N	O	S	0	0
			163	99	25	35	4		
2	I	21	Total	C	N	O	S	0	0
			163	99	25	35	4		

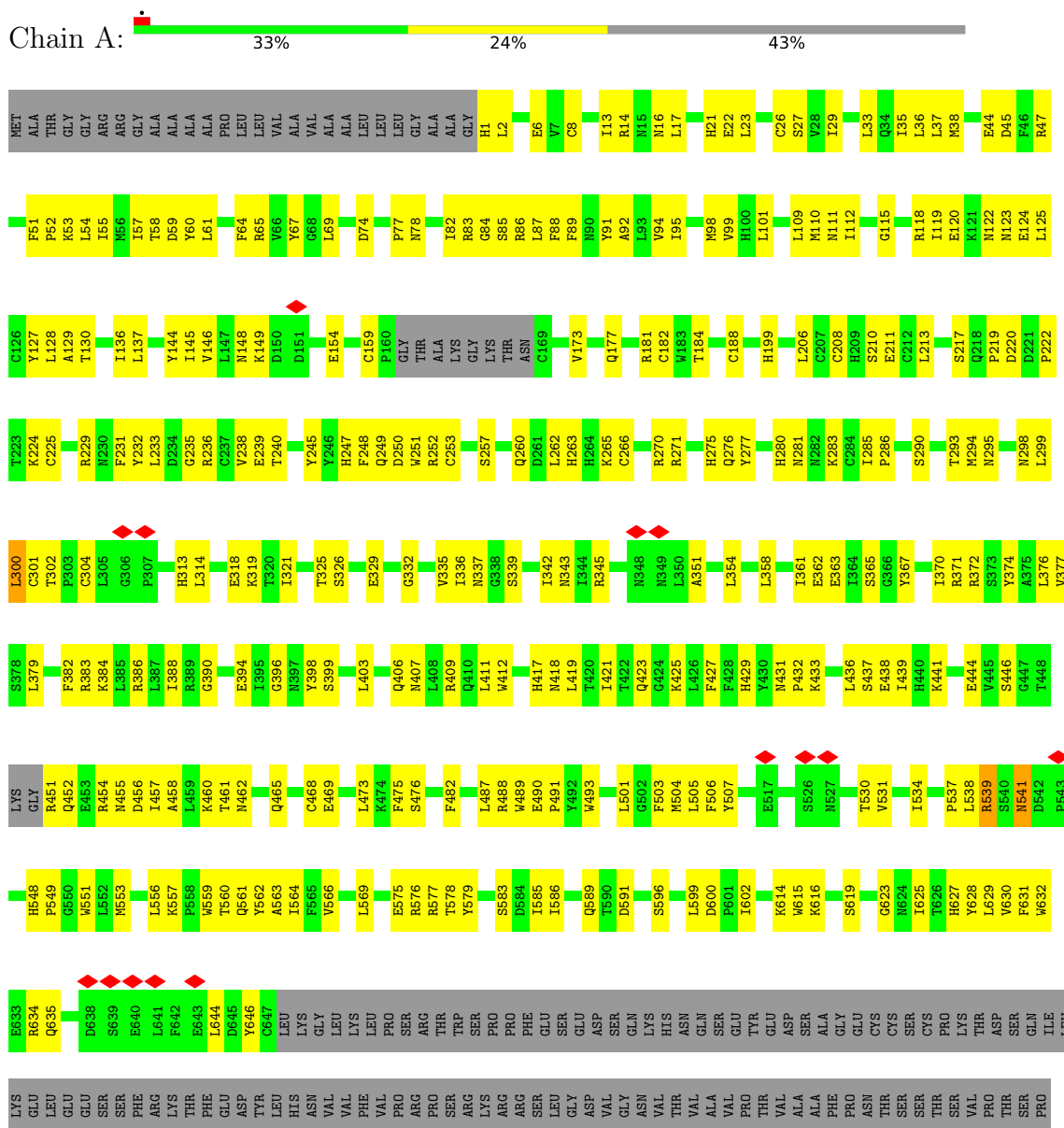
- Molecule 3 is a protein called Insulin.

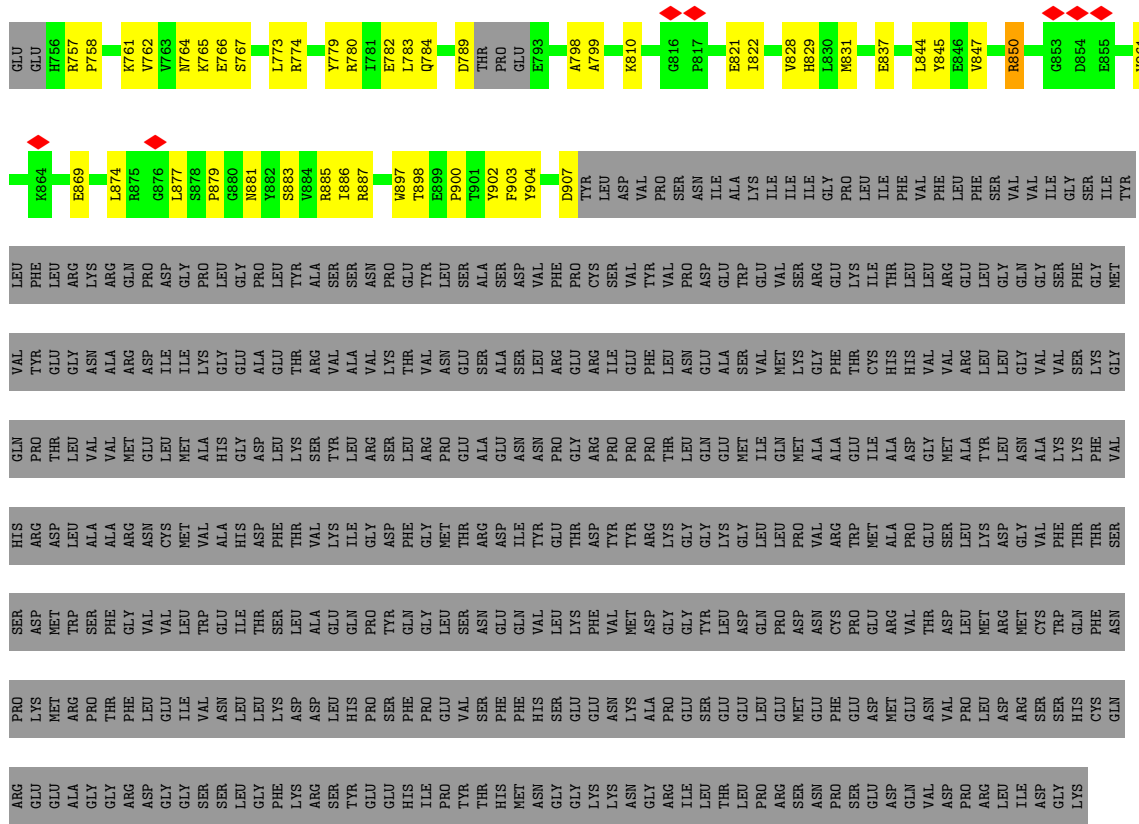
Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	25	Total	C	N	O	S	0	0
			200	129	34	35	2		
3	J	23	Total	C	N	O	S	0	0
			185	121	31	31	2		

3 Residue-property plots

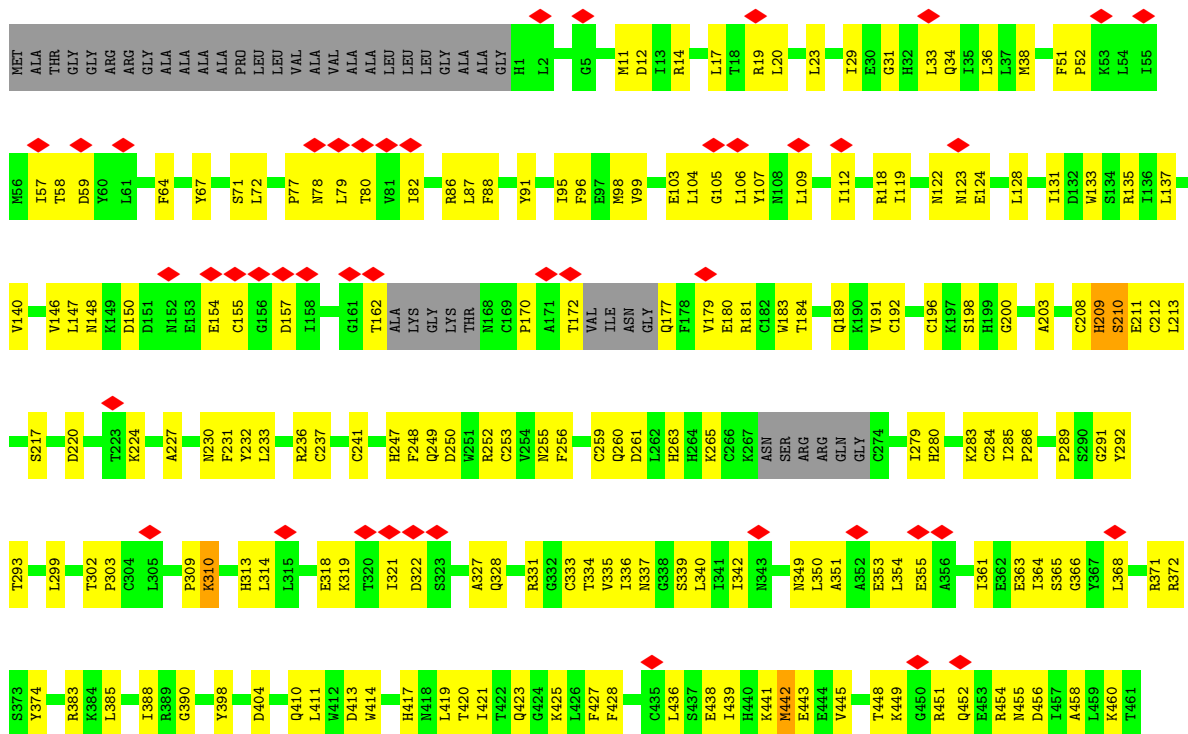
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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.

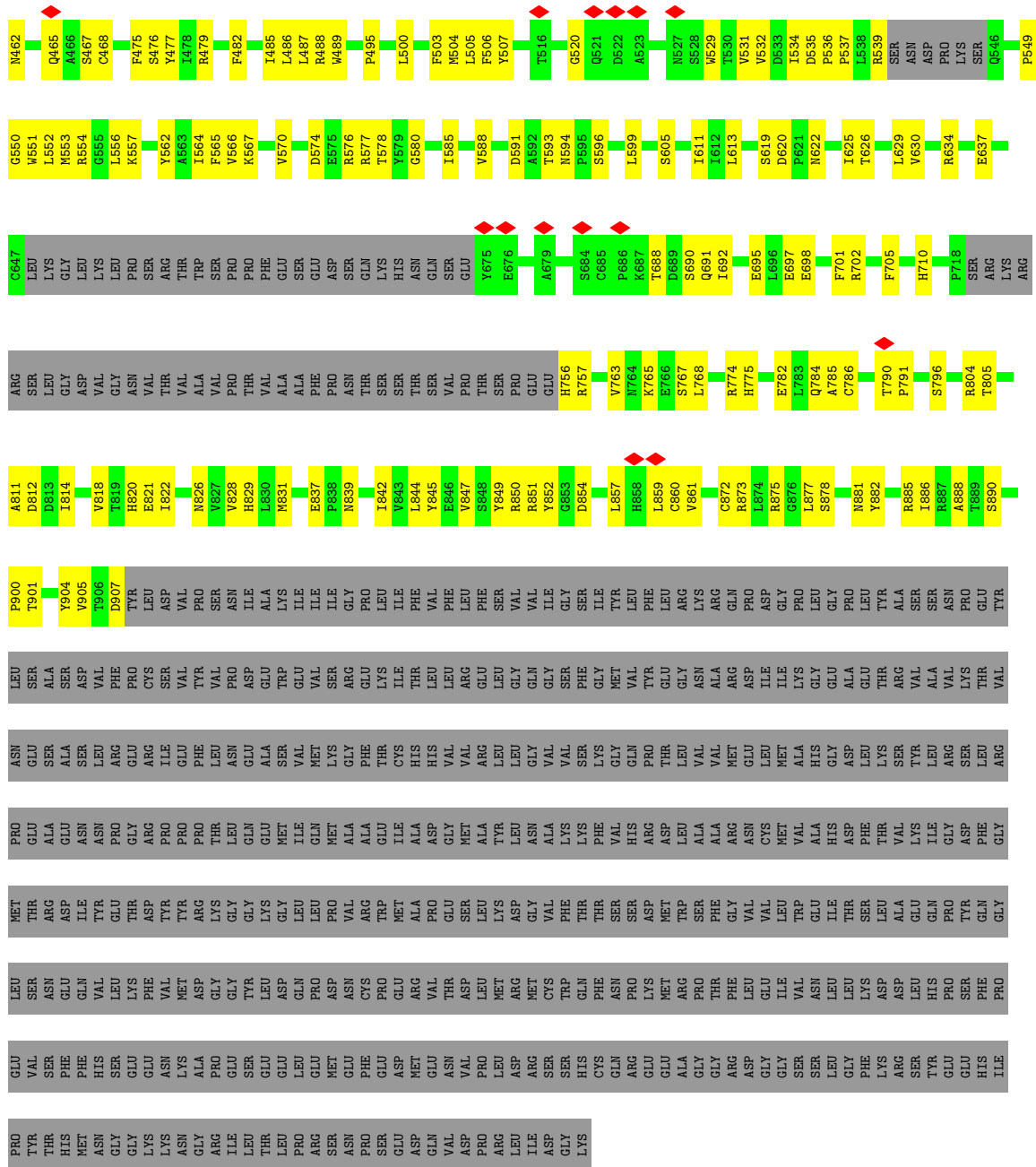
• Molecule 1: Isoform Short of Insulin receptor



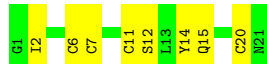


● Molecule 1: Isoform Short of Insulin receptor



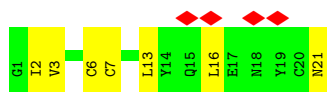


• Molecule 2: Insulin



• Molecule 2: Insulin





• Molecule 3: Insulin



• Molecule 3: Insulin



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	26701	Depositor
Resolution determination method	OTHER	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)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.400	Depositor
Minimum map value	-0.091	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.015	Depositor
Recommended contour level	0.125	Depositor
Map size (Å)	457.716, 457.716, 457.716	wwPDB
Map dimensions	420, 420, 420	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.0898, 1.0898, 1.0898	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.31	0/6496	0.68	2/8811 (0.0%)
1	B	0.28	0/6794	0.67	2/9213 (0.0%)
2	C	0.30	0/164	0.54	0/220
2	I	0.28	0/164	0.58	0/220
3	D	0.38	0/205	0.61	0/276
3	J	0.26	0/190	0.60	0/255
All	All	0.29	0/14013	0.67	4/18995 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	831	MET	CA-CB-CG	6.75	124.77	113.30
1	A	300	LEU	CA-CB-CG	6.01	129.12	115.30
1	B	442	MET	CA-CB-CG	5.16	122.07	113.30
1	B	504	MET	CA-CB-CG	5.15	122.06	113.30

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	A	6340	0	6134	232	0
1	B	6631	0	6398	252	0
2	C	163	0	149	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	163	0	149	9	0
3	D	200	0	184	14	0
3	J	185	0	171	11	0
All	All	13682	0	13185	497	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 19.

The worst 5 of 497 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:135:ARG:NH1	1:B:192:CYS:SG	2.37	0.97
1:A:95:ILE:HD11	1:A:119:ILE:HG23	1.59	0.83
1:A:437:SER:HA	1:A:577:ARG:HH22	1.44	0.81
1:A:336:ILE:HG12	1:A:361:ILE:HG21	1.63	0.81
1:A:266:CYS:HB2	1:A:271:ARG:HH21	1.47	0.79

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	776/1382 (56%)	763 (98%)	11 (1%)	2 (0%)	41	77
1	B	808/1382 (58%)	799 (99%)	6 (1%)	3 (0%)	34	72
2	C	19/21 (90%)	19 (100%)	0	0	100	100
2	I	19/21 (90%)	17 (90%)	2 (10%)	0	100	100
3	D	23/30 (77%)	23 (100%)	0	0	100	100
3	J	21/30 (70%)	21 (100%)	0	0	100	100
All	All	1666/2866 (58%)	1642 (99%)	19 (1%)	5 (0%)	44	77

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	54	LEU
1	B	208	CYS
1	B	210	SER
1	A	541	ASN
1	B	310	LYS

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	713/1226 (58%)	706 (99%)	7 (1%)	76 86
1	B	747/1226 (61%)	744 (100%)	3 (0%)	91 94
2	C	20/20 (100%)	20 (100%)	0	100 100
2	I	20/20 (100%)	20 (100%)	0	100 100
3	D	21/26 (81%)	21 (100%)	0	100 100
3	J	19/26 (73%)	19 (100%)	0	100 100
All	All	1540/2544 (60%)	1530 (99%)	10 (1%)	86 92

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	209	HIS
1	B	210	SER
1	B	211	GLU
1	A	596	SER
1	A	625	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	622	ASN
1	B	775	HIS

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Mol	Chain	Res	Type
1	B	839	ASN
1	A	775	HIS
1	A	784	GLN

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.

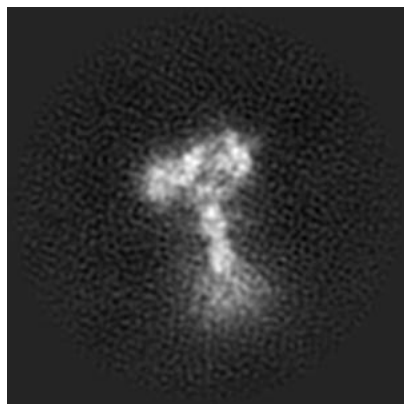
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13388. These allow visual inspection of the internal detail of the map and identification of artifacts.

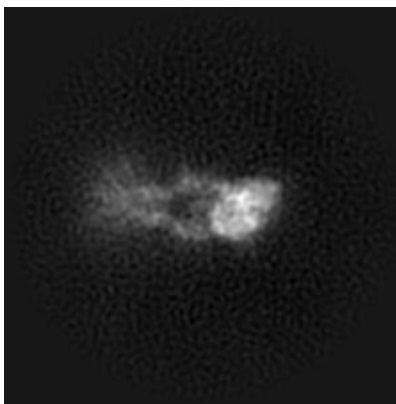
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

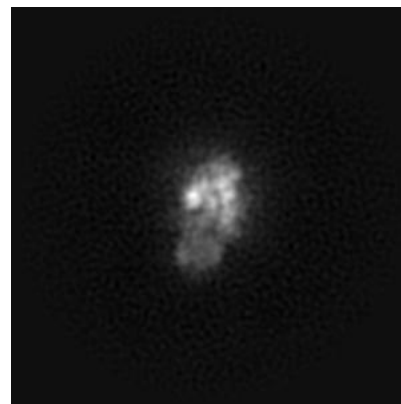
6.1.1 Primary map



X

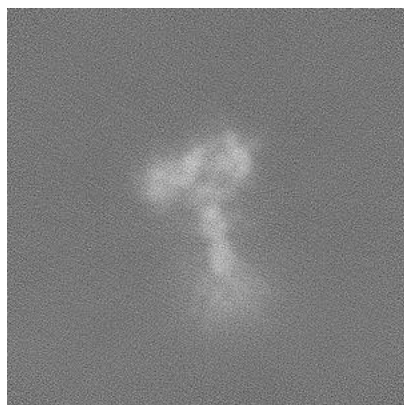


Y

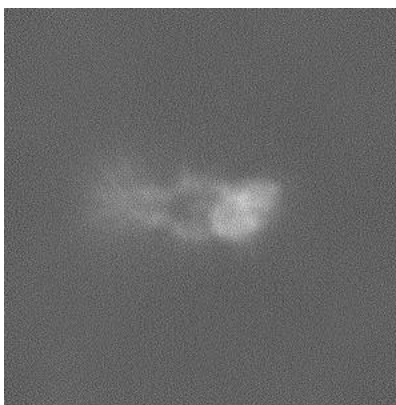


Z

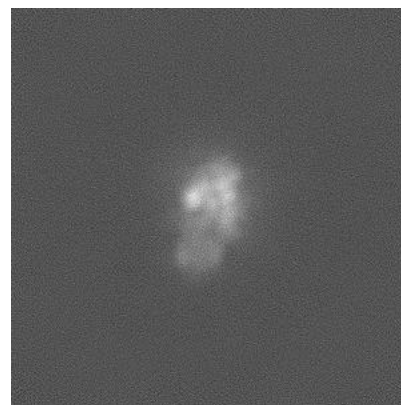
6.1.2 Raw map



X



Y

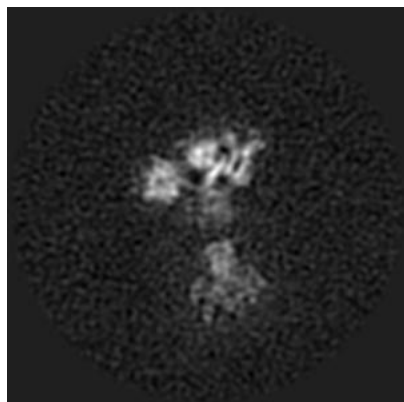


Z

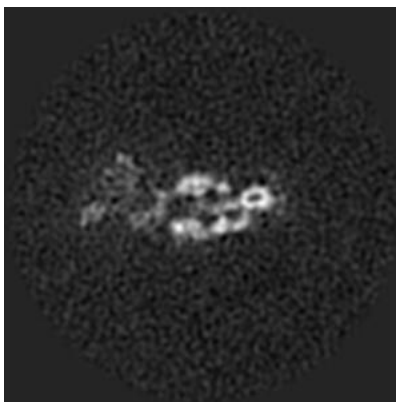
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

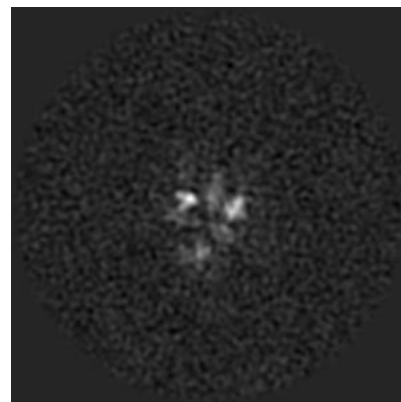
6.2.1 Primary map



X Index: 210

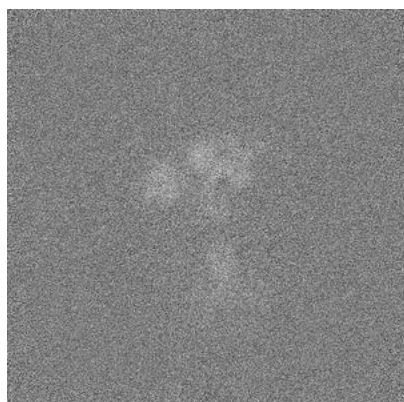


Y Index: 210

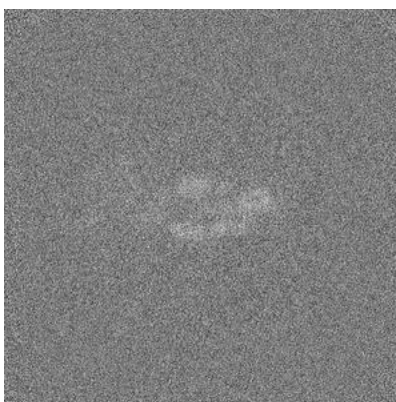


Z Index: 210

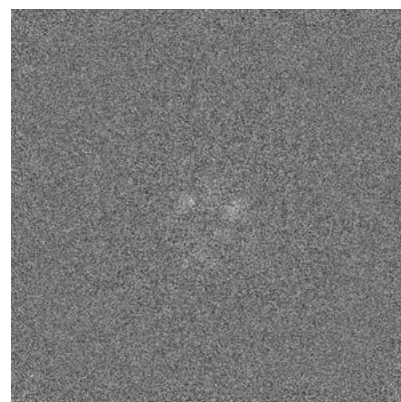
6.2.2 Raw map



X Index: 210



Y Index: 210

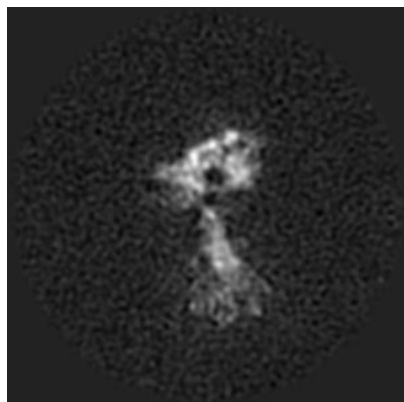


Z Index: 210

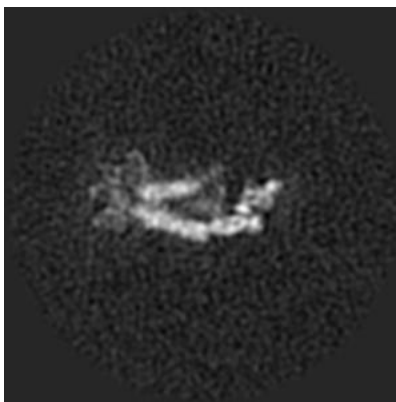
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

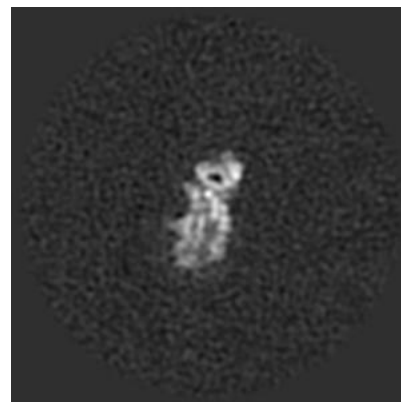
6.3.1 Primary map



X Index: 223

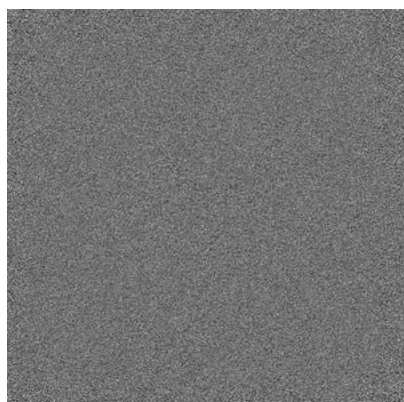


Y Index: 221

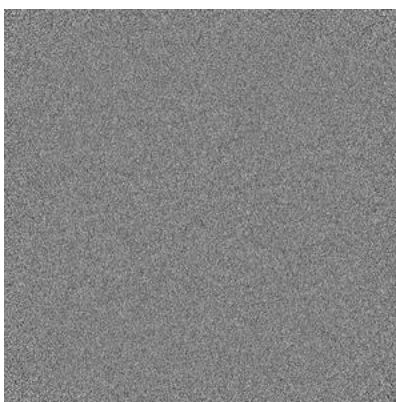


Z Index: 253

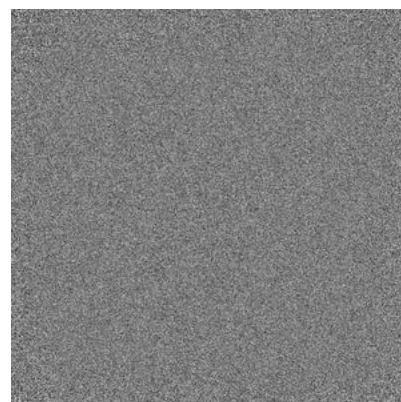
6.3.2 Raw map



X Index: 0



Y Index: 0



Z Index: 0

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

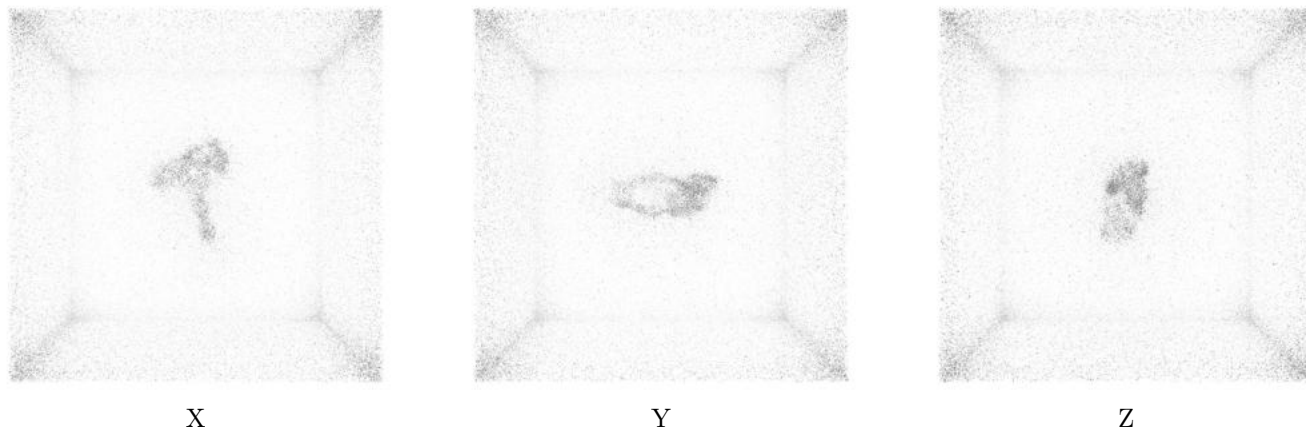
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.125. 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



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.

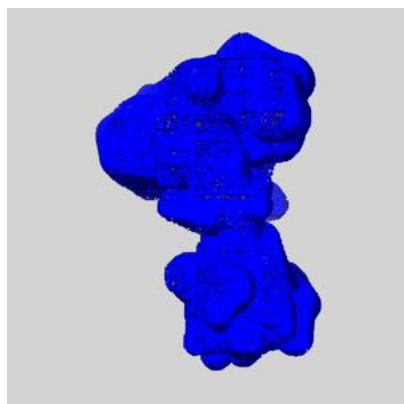
6.5 Mask visualisation [i](#)

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

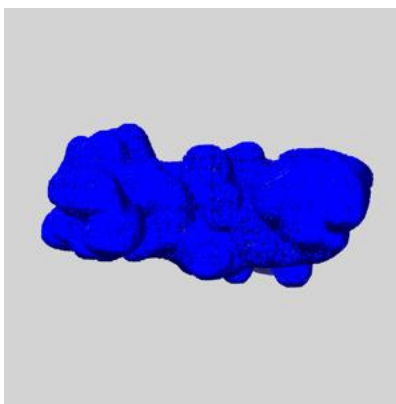
A mask typically either:

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

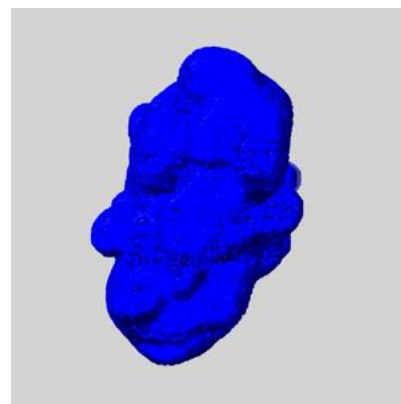
6.5.1 emd_13388_msk_1.map [i](#)



X



Y

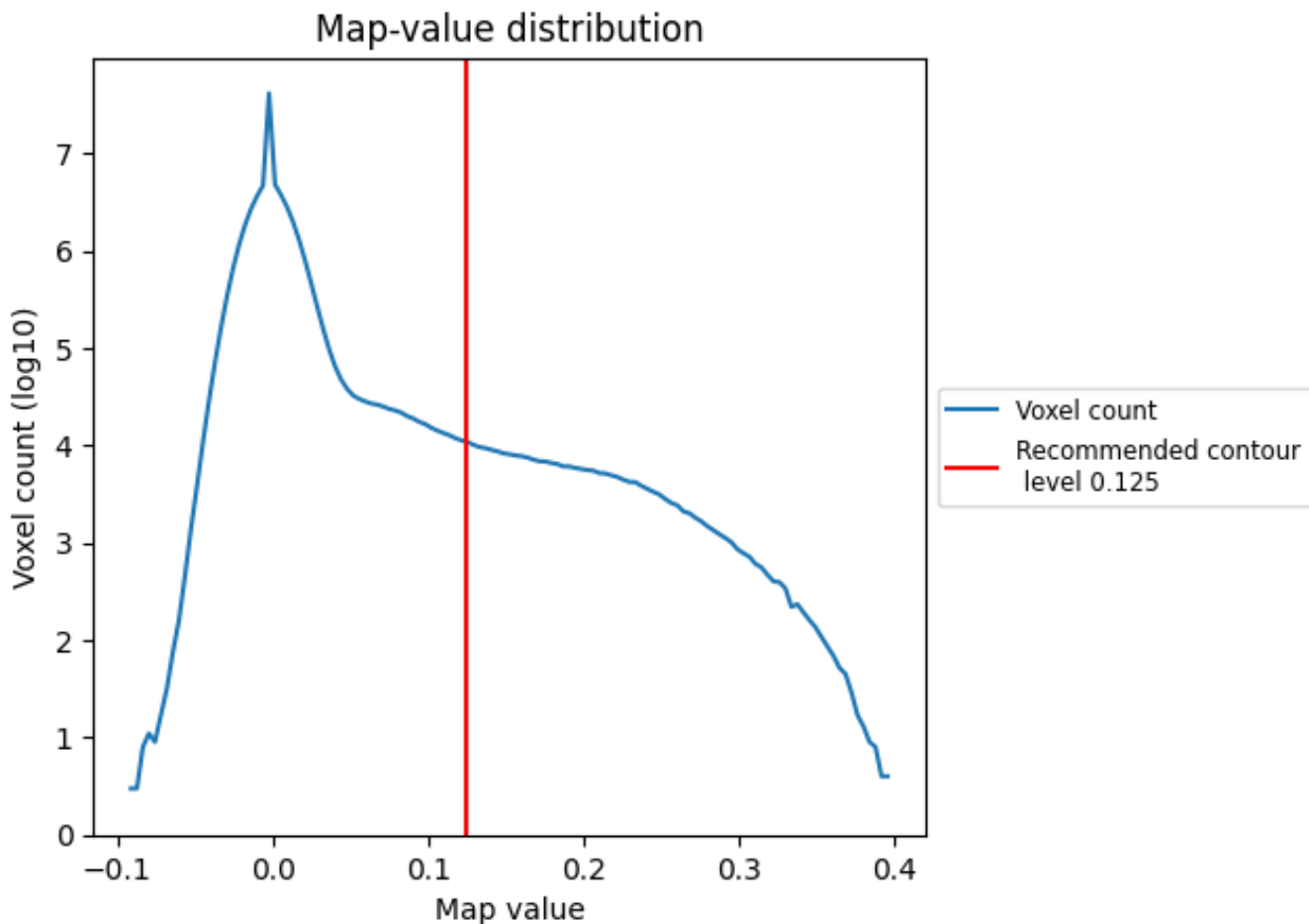


Z

7 Map analysis [i](#)

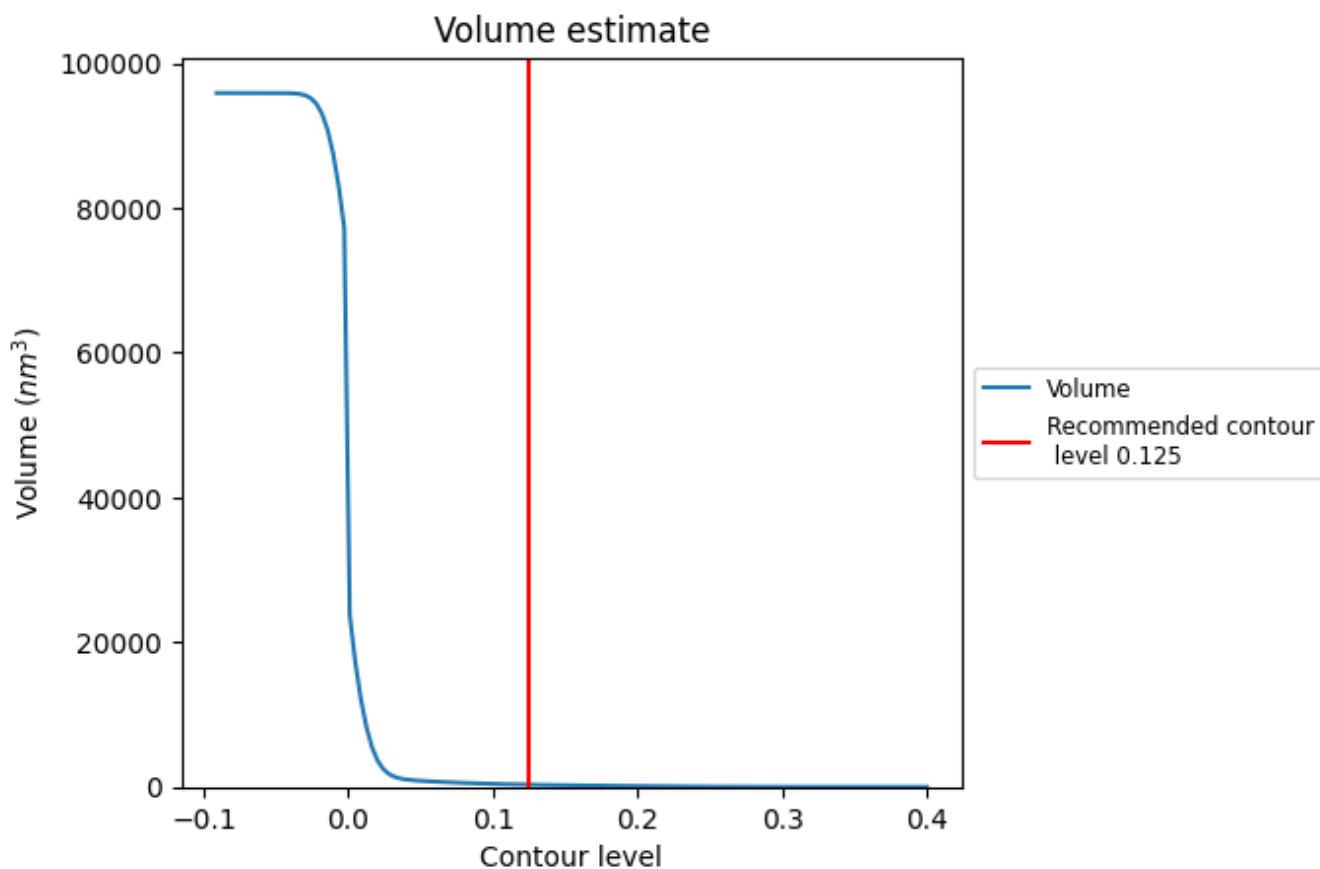
This section contains the results of statistical analysis of the map.

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.

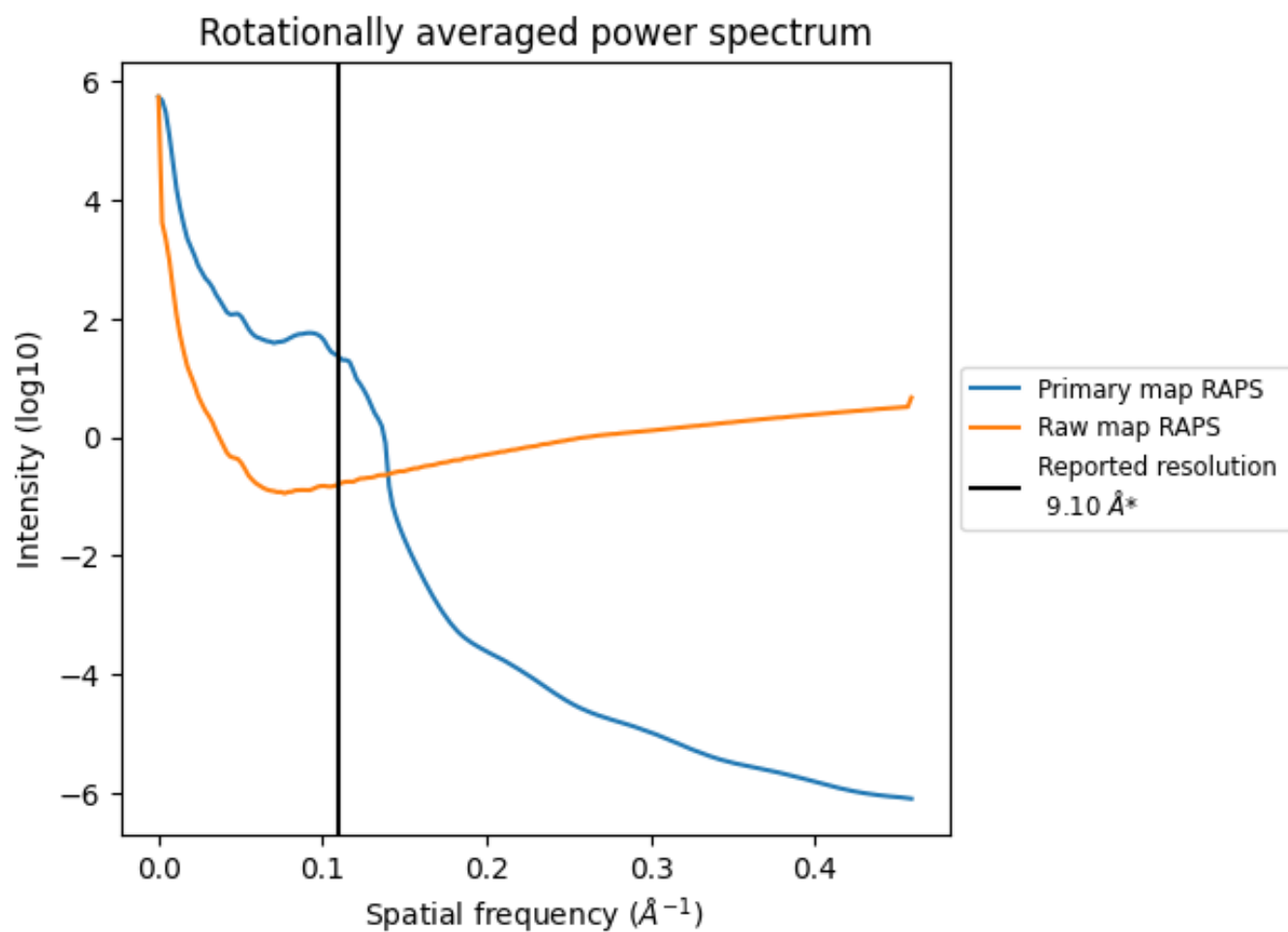
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 314 nm^3 ; this corresponds to an approximate mass of 284 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](#)

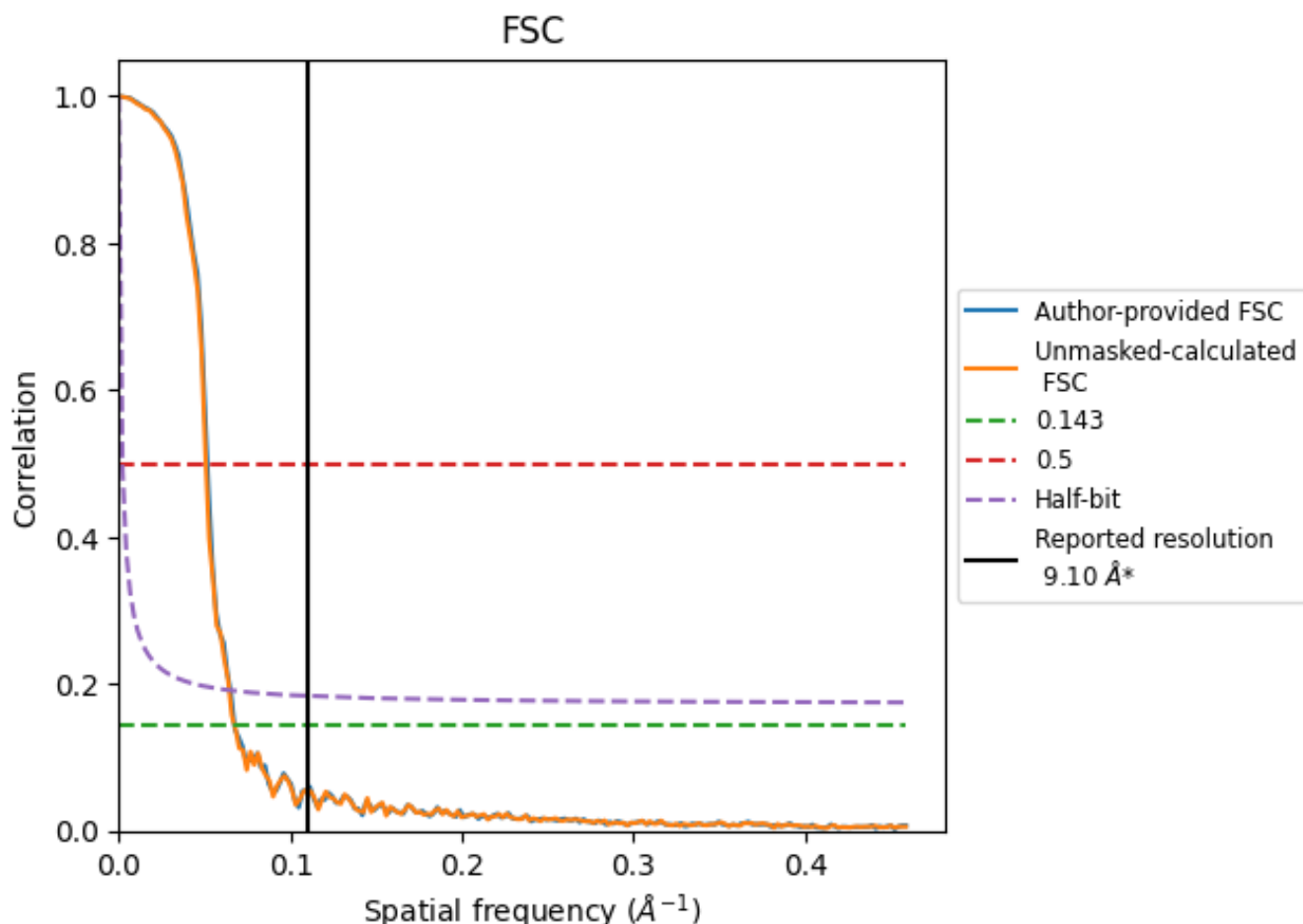


*Reported resolution corresponds to spatial frequency of 0.110 Å⁻¹

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 0.110 Å⁻¹

8.2 Resolution estimates [i](#)

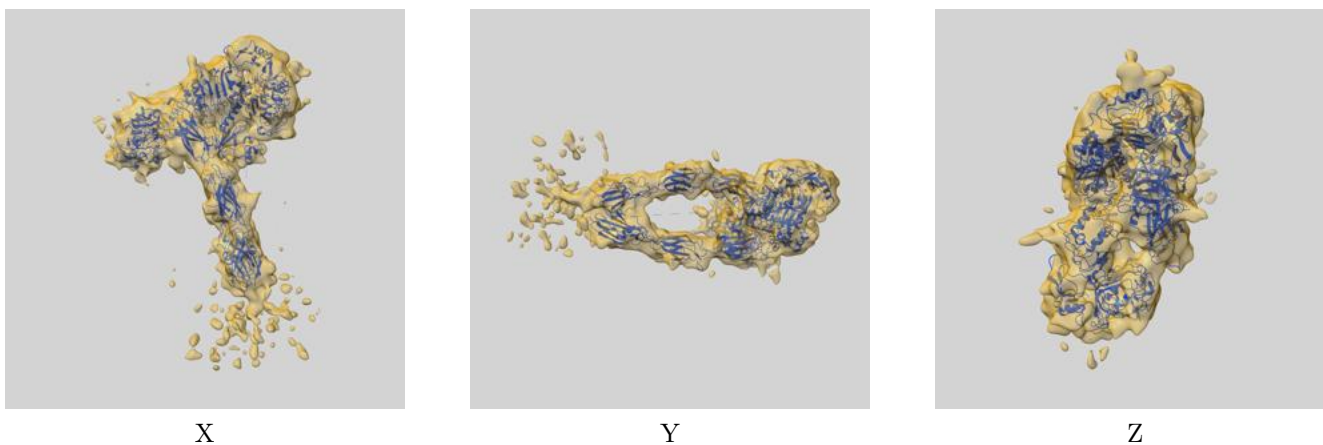
Resolution estimate (Å)	Estimation criterion (FSC cut-off)			
	0.143	0.5	Half-bit	Other
Reported by author	-	-	-	9.10
Author-provided FSC curve	14.81	19.34	15.36	-
Unmasked-calculated*	14.84	19.69	15.55	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

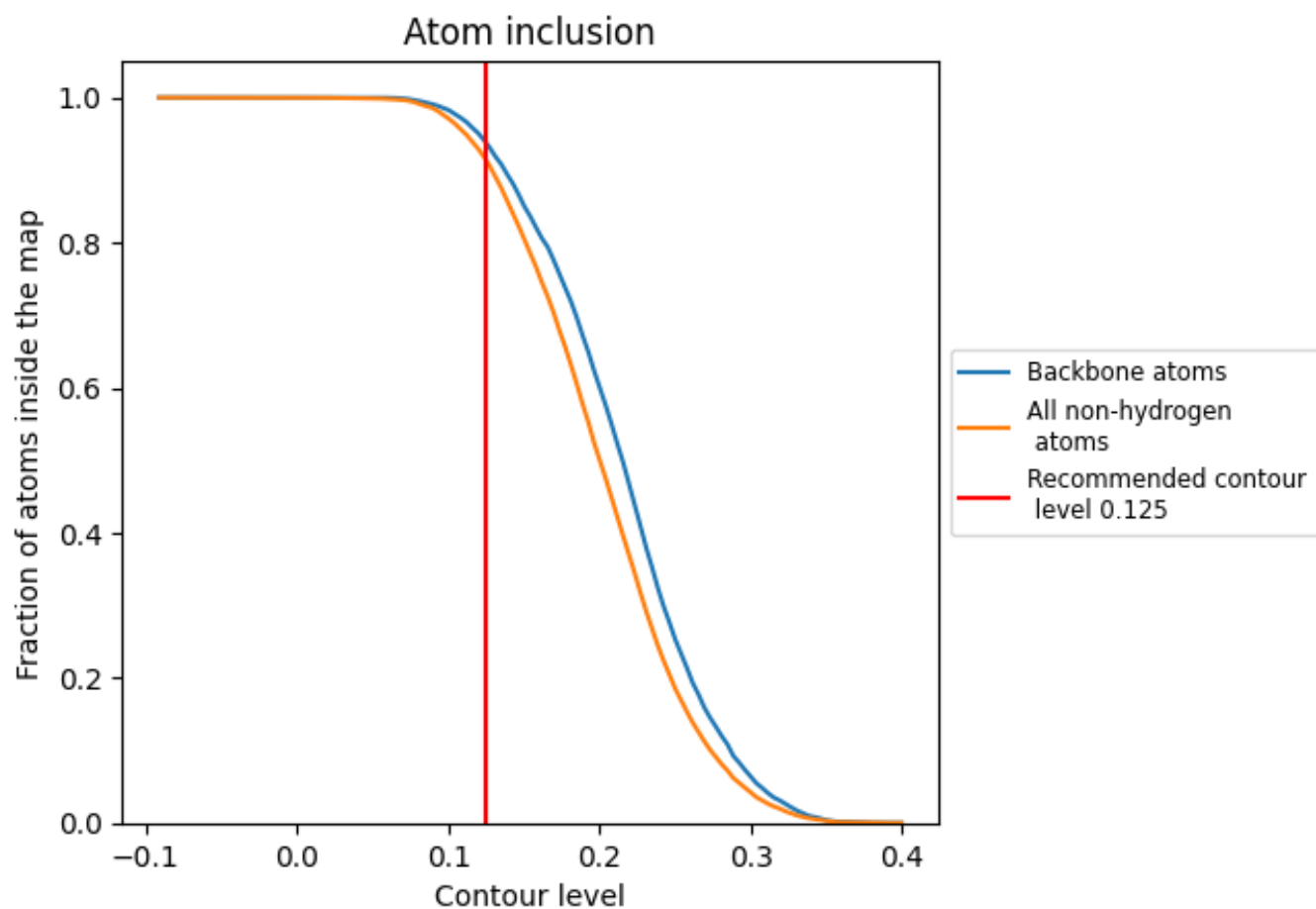
This section contains information regarding the fit between EMDB map EMD-13388 and PDB model 7PG4. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.125 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 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 91% of all non-hydrogen atoms, are inside the map.