



wwPDB EM Validation Summary Report ⓘ

Jan 9, 2024 – 04:27 PM JST

PDB ID : 8JA3
EMDB ID : EMD-36124
Title : Structure of beta-arrestin1 in complex with C3aRpp
Authors : Maharana, J.; Sarma, P.; Yadav, M.K.; Chami, M.; Banerjee, R.; Shukla, A.K.
Deposited on : 2023-05-05
Resolution : 3.94 Å (reported)
Based on initial model : 8GO8

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 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:

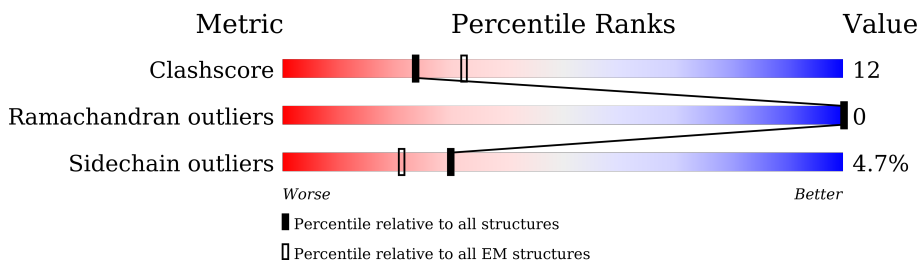
EMDB validation analysis : 0.0.1.dev70
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance

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

The reported resolution of this entry is 3.94 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	357	<p>65% 20% 13%</p>
1	B	357	<p>70% 17% 11%</p>
2	U	10	<p>60% 30% 10%</p>
2	V	10	<p>60% 30% 10%</p>
3	H	237	<p>32% 10% 58%</p>
3	I	237	<p>34% 10% 55%</p>
4	L	215	<p>40% 6% 53%</p>
4	M	215	<p>43% 6% 53%</p>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7130 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 Beta-arrestin-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	B	316	Total	C	N	O	S	0	0
			2118	1363	367	382	6		
1	A	309	Total	C	N	O	S	0	0
			2070	1323	357	384	6		

- Molecule 2 is a protein called C3a anaphylatoxin chemotactic receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	N	O	P			S
2	U	10	Total	C	N	O	P	S	0	0
			81	38	12	26	4	1		
2	V	10	Total	C	N	O	P	S	0	0
			81	38	12	26	4	1		

- Molecule 3 is a protein called Fab30 heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	I	107	Total	C	N	O	S	0	0
			756	483	128	142	3		
3	H	99	Total	C	N	O	S	0	0
			684	438	119	125	2		

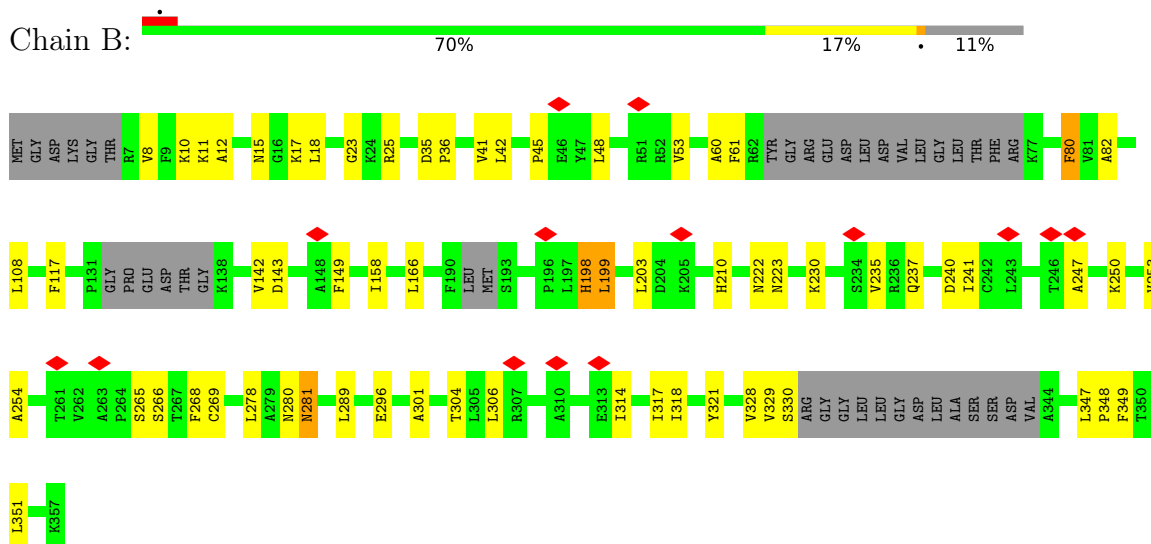
- Molecule 4 is a protein called Fab30 light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	M	102	Total	C	N	O	S	0	0
			680	436	115	127	2		
4	L	100	Total	C	N	O	S	0	0
			660	418	112	128	2		

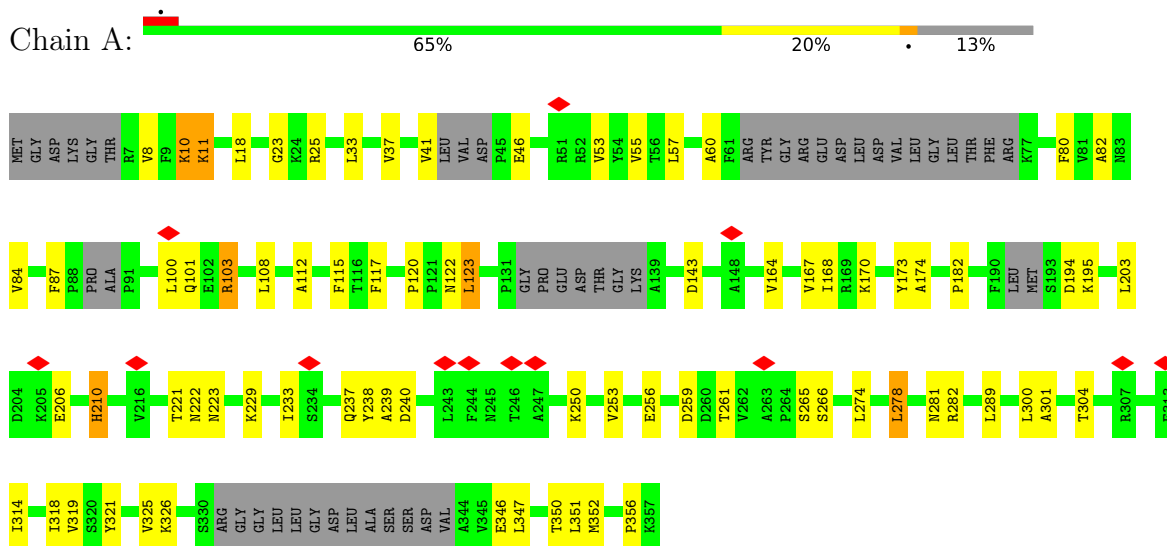
3 Residue-property plots

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: Beta-arrestin-1



- Molecule 1: Beta-arrestin-1



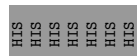
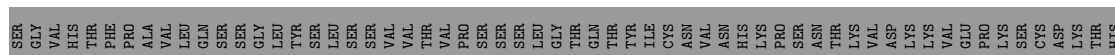
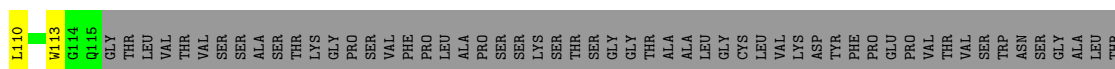
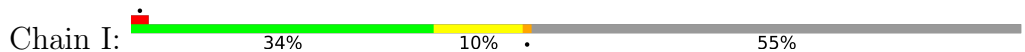
- Molecule 2: C3a anaphylatoxin chemotactic receptor



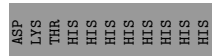
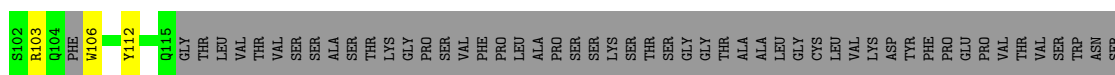
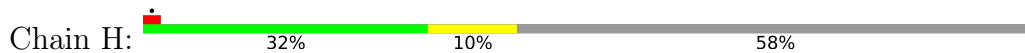
- Molecule 2: C3a anaphylatoxin chemotactic receptor



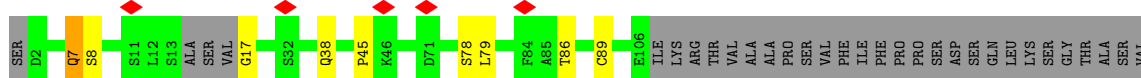
- Molecule 3: Fab30 heavy chain



- Molecule 3: Fab30 heavy chain



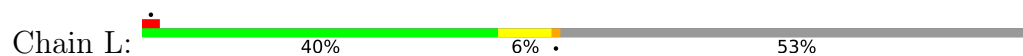
- Molecule 4: Fab30 light chain



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

ALA
CYS
GLU
VAL
THR
HIS
GLN
LEU
SER
PRO
VAL
THR
LYS
SER
PHE
ASN
ARG
GLY
GLU
CYS

● Molecule 4: Fab30 light chain



SER
D2
L12
S13
ALA
SER
VAL
G17
S29
Y34
A35
W36
Y37
K40
F41
K46
S57
D71
F72
L79
PRO
E82
D83
F84
Y87
Y88
V97
T98
E106
ILE
LYS
SER
THR
SER
VAL
THR
ALA
ALA
PRO
SER
VAL
PHE
ILE
PHE
PRO
LEU
SER
ASP
SER
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CYS
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ASN
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ARG
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TRP
LYS
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VAL
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THR
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GLY
LEU
SER
SER
PRO
VAL
THR
LYS
SER
PHE
ASN
ARG
GLY
GLU
CYS

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	164339	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS GLACIOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	49.4	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	46000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	2.037	Depositor
Minimum map value	-0.034	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.013	Depositor
Recommended contour level	0.044	Depositor
Map size (\AA)	365.2608, 365.2608, 365.2608	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.4268, 1.4268, 1.4268	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: TPO, 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	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.35	1/2109 (0.0%)	0.52	0/2902
1	B	0.32	0/2161	0.52	0/2979
2	U	0.33	0/37	1.05	0/46
2	V	0.37	0/37	0.53	0/46
3	H	0.35	0/700	0.53	0/958
3	I	0.35	0/774	0.55	0/1057
4	L	0.34	0/672	0.57	0/922
4	M	0.31	0/696	0.52	0/958
All	All	0.34	1/7186 (0.0%)	0.53	0/9868

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	33	LEU	C-N	5.83	1.47	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2070	0	1778	55	0
1	B	2118	0	1845	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	U	81	0	48	5	0
2	V	81	0	47	3	0
3	H	684	0	558	14	0
3	I	756	0	653	32	0
4	L	660	0	531	10	0
4	M	680	0	556	5	0
All	All	7130	0	6016	159	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 12.

The worst 5 of 159 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)
3:I:57:TYR:HB3	3:I:103:ARG:HH22	1.44	0.81
3:I:57:TYR:CB	3:I:103:ARG:NH2	2.47	0.77
3:I:110:LEU:HD22	3:I:113:TRP:CE2	2.20	0.76
3:I:100:ALA:HB1	3:I:110:LEU:HD21	1.69	0.73
3:I:57:TYR:HB3	3:I:103:ARG:NH2	2.03	0.73

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	295/357 (83%)	272 (92%)	23 (8%)	0	100	100
1	B	306/357 (86%)	285 (93%)	21 (7%)	0	100	100
2	U	5/10 (50%)	3 (60%)	2 (40%)	0	100	100
2	V	5/10 (50%)	4 (80%)	1 (20%)	0	100	100
3	H	91/237 (38%)	83 (91%)	8 (9%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	I	101/237 (43%)	95 (94%)	6 (6%)	0	100	100
4	L	94/215 (44%)	85 (90%)	9 (10%)	0	100	100
4	M	98/215 (46%)	93 (95%)	5 (5%)	0	100	100
All	All	995/1638 (61%)	920 (92%)	75 (8%)	0	100	100

There are no Ramachandran outliers to report.

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	173/315 (55%)	163 (94%)	10 (6%)	20	48
1	B	176/315 (56%)	170 (97%)	6 (3%)	37	61
2	U	3/6 (50%)	3 (100%)	0	100	100
2	V	3/6 (50%)	3 (100%)	0	100	100
3	H	52/200 (26%)	51 (98%)	1 (2%)	57	75
3	I	65/200 (32%)	61 (94%)	4 (6%)	18	46
4	L	51/190 (27%)	48 (94%)	3 (6%)	19	47
4	M	53/190 (28%)	50 (94%)	3 (6%)	20	49
All	All	576/1422 (40%)	549 (95%)	27 (5%)	30	53

5 of 27 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LYS
1	A	103	ARG
4	L	12	LEU
1	A	84	VAL
1	A	123	LEU

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

Mol	Chain	Res	Type
1	A	237	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](#)

8 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
2	TPO	V	466	2	8,10,11	0.69	0	10,14,16	1.27	1 (10%)
2	SEP	V	465	2	8,9,10	0.61	0	8,12,14	0.64	0
2	SEP	U	459	2	8,9,10	1.55	1 (12%)	8,12,14	1.35	2 (25%)
2	SEP	V	459	2	8,9,10	1.55	1 (12%)	8,12,14	1.35	2 (25%)
2	TPO	U	463	2	8,10,11	0.60	0	10,14,16	0.96	1 (10%)
2	TPO	V	463	2	8,10,11	0.63	0	10,14,16	1.00	1 (10%)
2	SEP	U	465	2	8,9,10	0.63	0	8,12,14	0.77	0
2	TPO	U	466	2	8,10,11	1.55	1 (12%)	10,14,16	1.83	1 (10%)

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
2	TPO	V	466	2	-	5/9/11/13	-
2	SEP	V	465	2	-	2/5/8/10	-
2	SEP	U	459	2	-	1/5/8/10	-
2	SEP	V	459	2	-	1/5/8/10	-
2	TPO	U	463	2	-	2/9/11/13	-
2	TPO	V	463	2	-	0/9/11/13	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SEP	U	465	2	-	4/5/8/10	-
2	TPO	U	466	2	-	2/9/11/13	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	U	459	SEP	P-O1P	3.35	1.61	1.50
2	V	459	SEP	P-O1P	3.34	1.61	1.50
2	U	466	TPO	P-O1P	3.21	1.60	1.50

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	U	466	TPO	P-OG1-CB	-5.11	107.77	123.21
2	V	466	TPO	O-C-CA	-3.10	116.65	124.78
2	V	459	SEP	OG-CB-CA	2.41	110.50	108.14
2	U	459	SEP	OG-CB-CA	2.40	110.48	108.14
2	V	463	TPO	O-C-CA	-2.34	118.64	124.78

There are no chirality outliers.

5 of 17 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	U	463	TPO	CB-OG1-P-O1P
2	U	465	SEP	N-CA-CB-OG
2	U	465	SEP	CB-OG-P-O1P
2	U	465	SEP	CB-OG-P-O2P
2	U	465	SEP	CB-OG-P-O3P

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	V	466	TPO	2	0
2	U	463	TPO	3	0

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-36124. These allow visual inspection of the internal detail of the map and identification of artifacts.

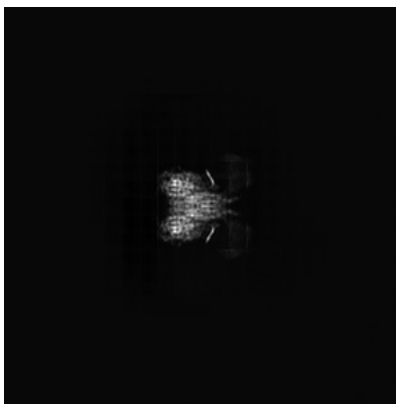
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

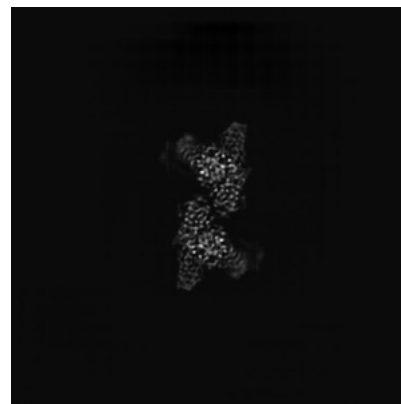
6.1.1 Primary map



X

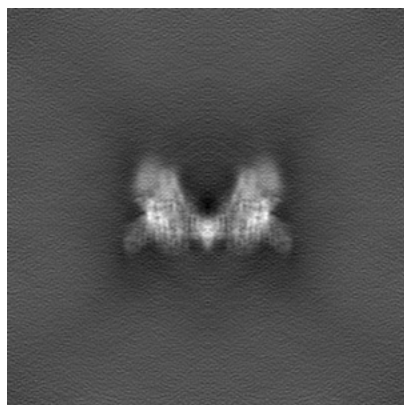


Y

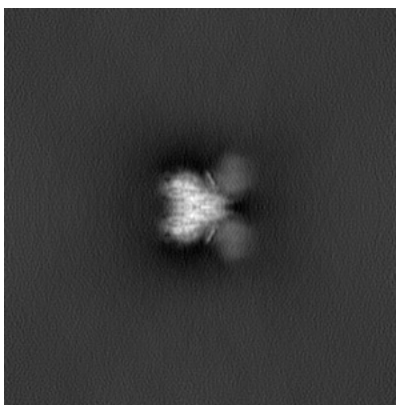


Z

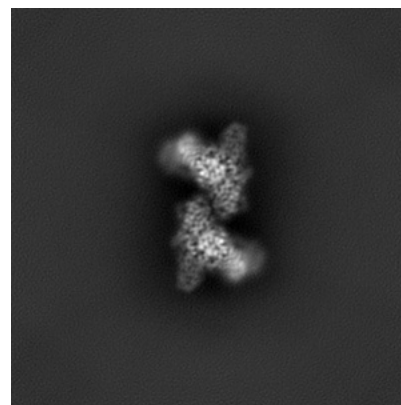
6.1.2 Raw map



X



Y

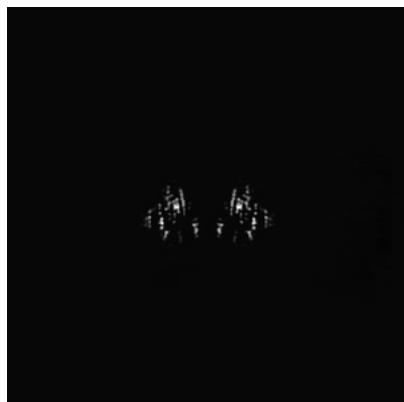


Z

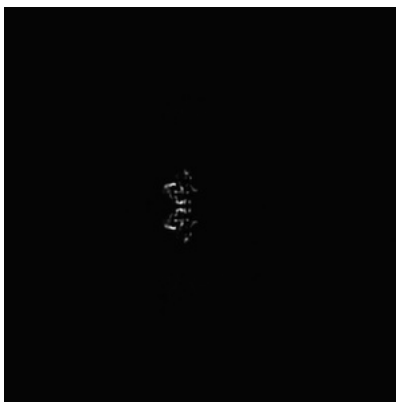
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 128

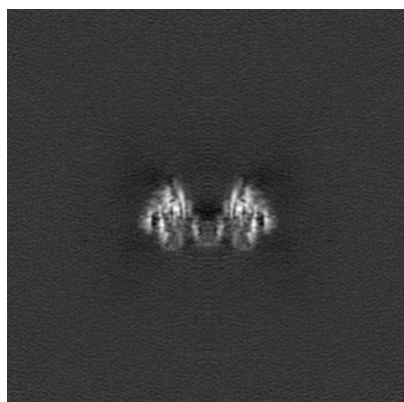


Y Index: 128

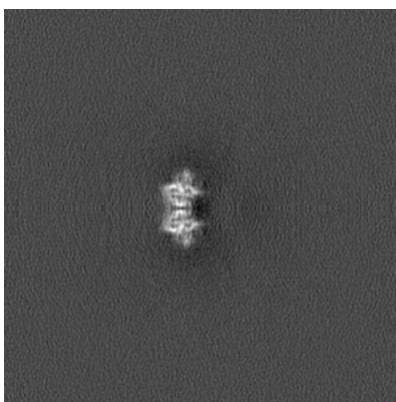


Z Index: 128

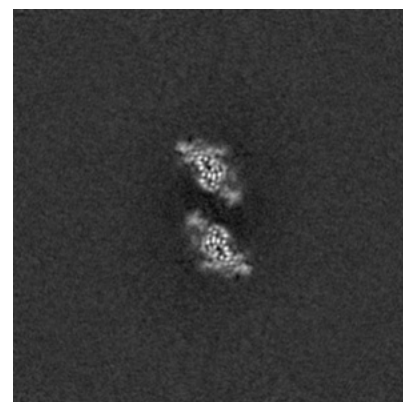
6.2.2 Raw map



X Index: 128



Y Index: 128



Z Index: 128

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

6.3 Largest variance slices [i](#)

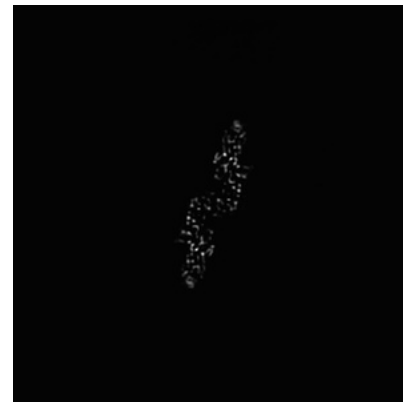
6.3.1 Primary map



X Index: 131

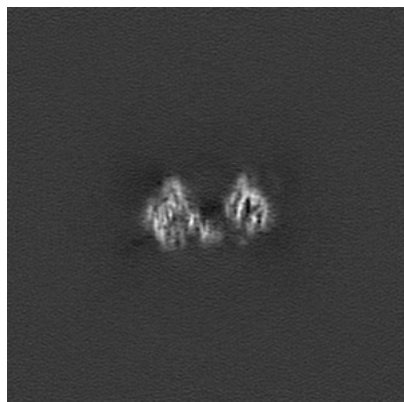


Y Index: 102

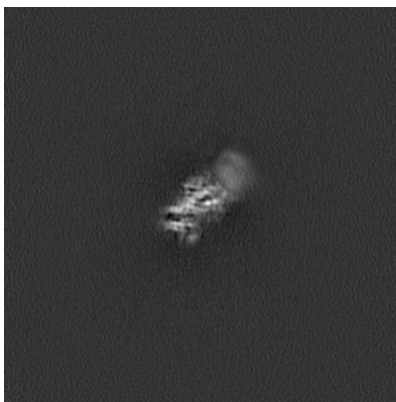


Z Index: 109

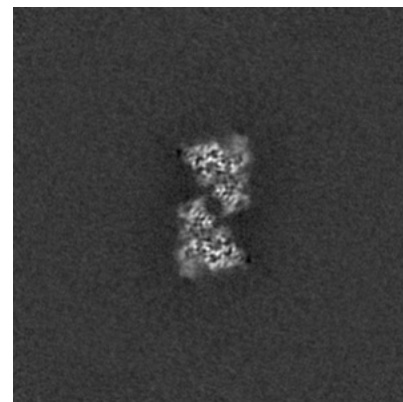
6.3.2 Raw map



X Index: 125



Y Index: 99

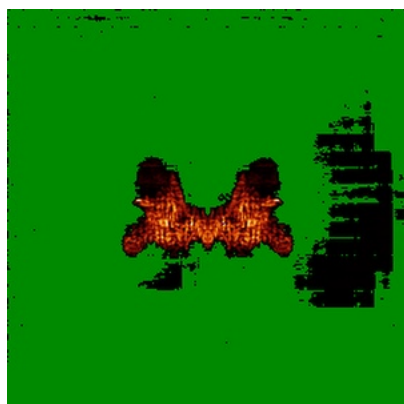


Z Index: 120

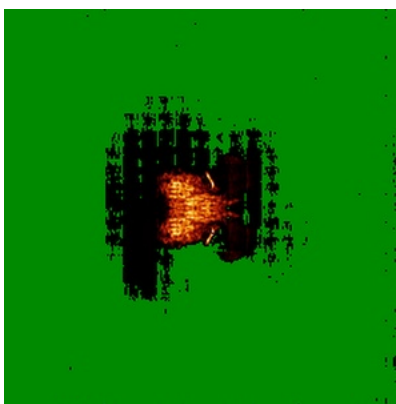
The images above show the largest variance slices of the map in three orthogonal directions.

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

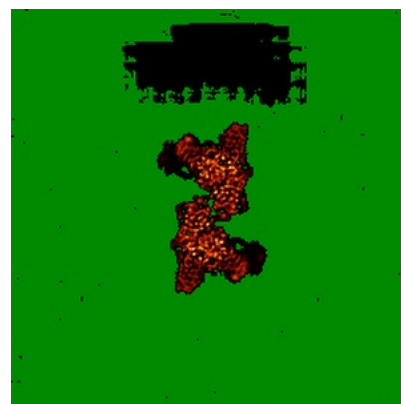
6.4.1 Primary map



X

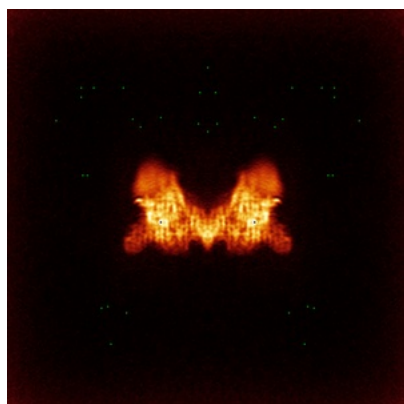


Y

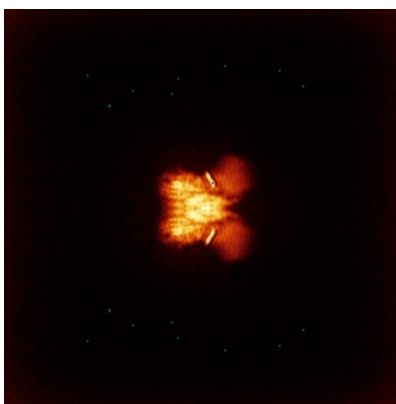


Z

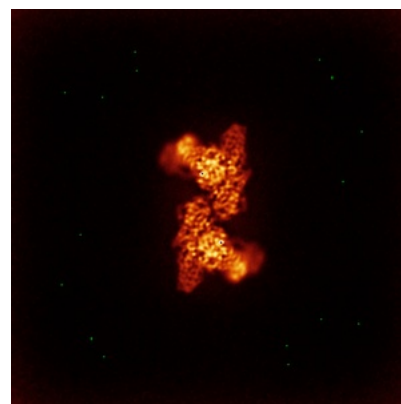
6.4.2 Raw map



X



Y

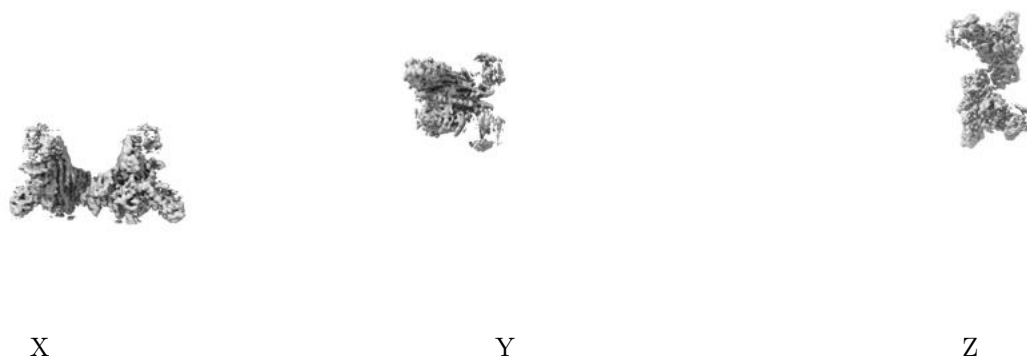


Z

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.

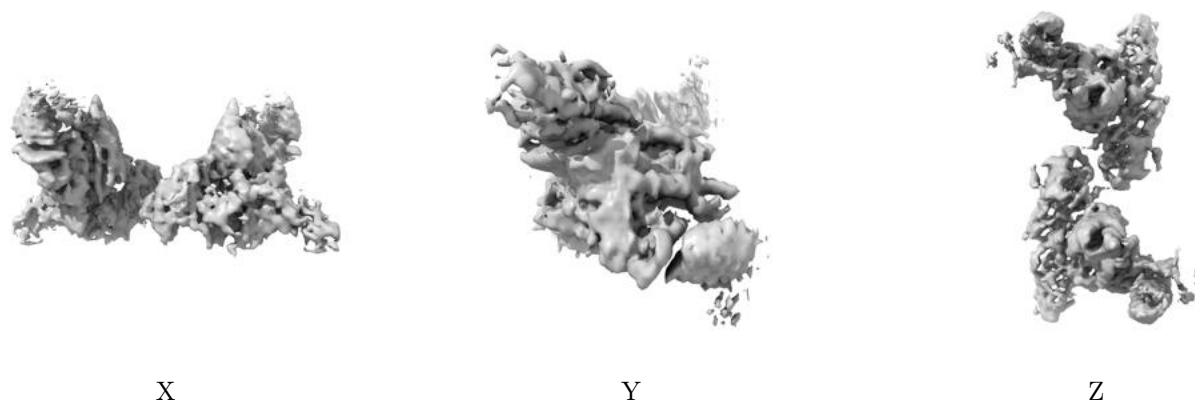
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.044. 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.

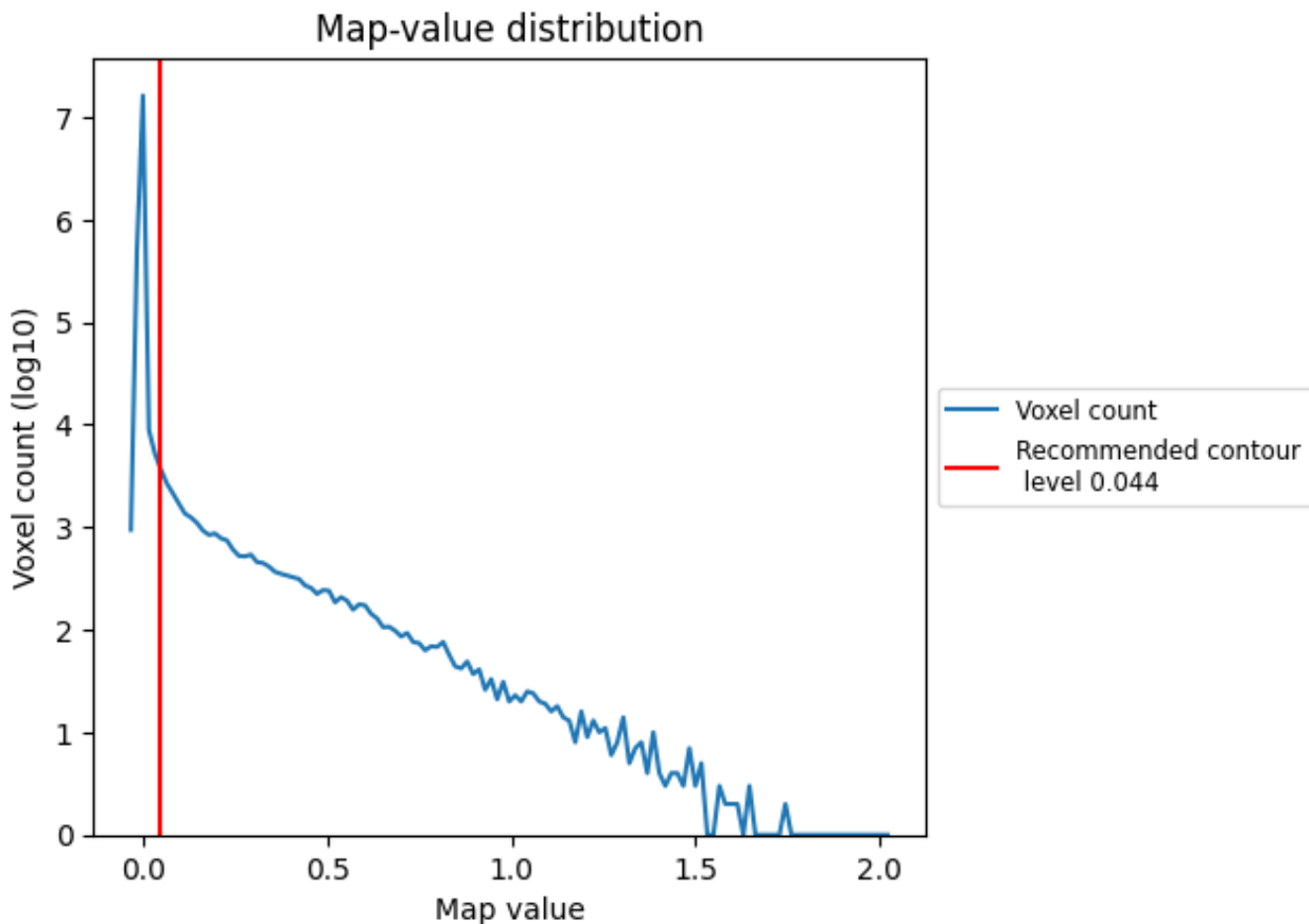
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

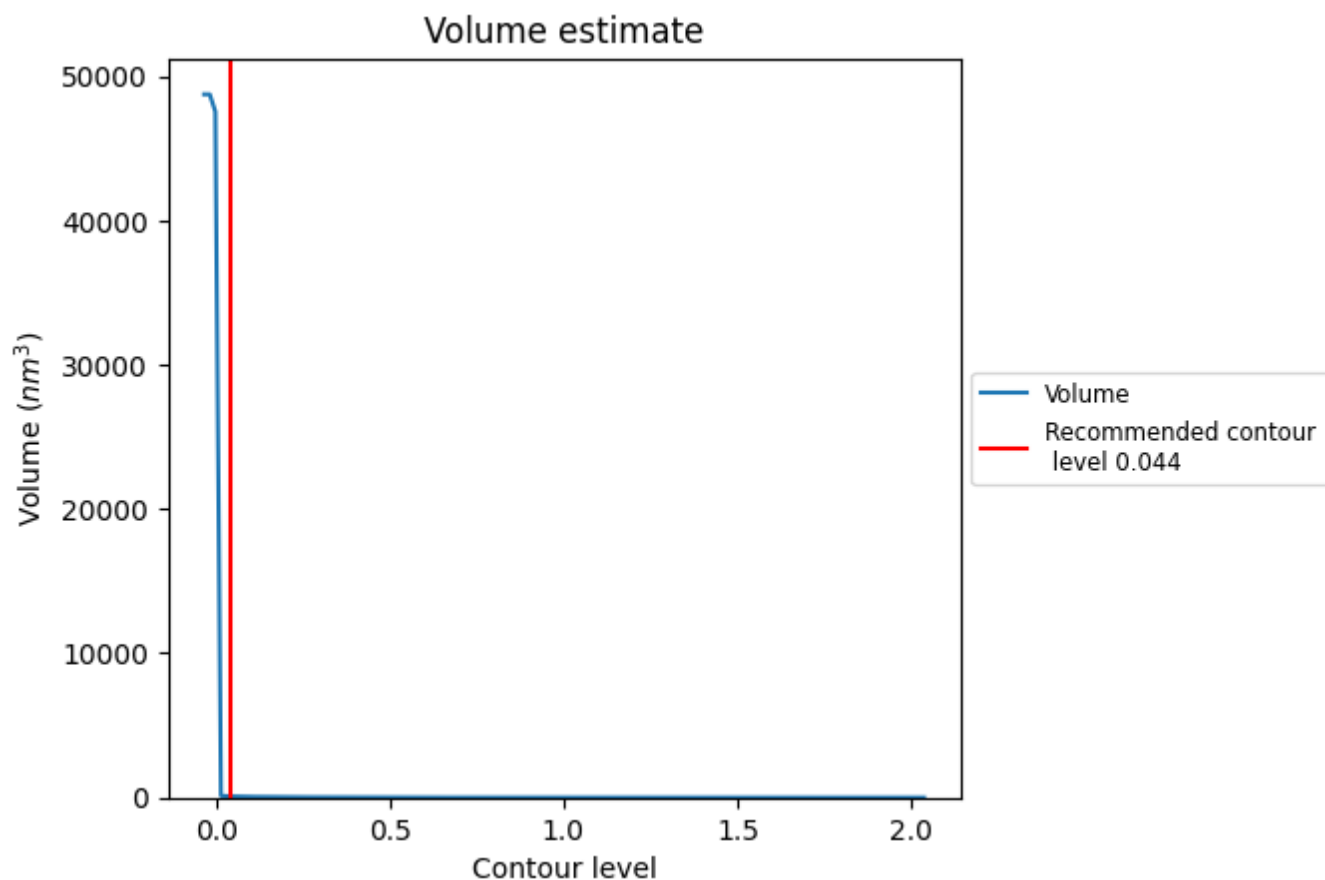
This section contains the results of statistical analysis of the map.

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.

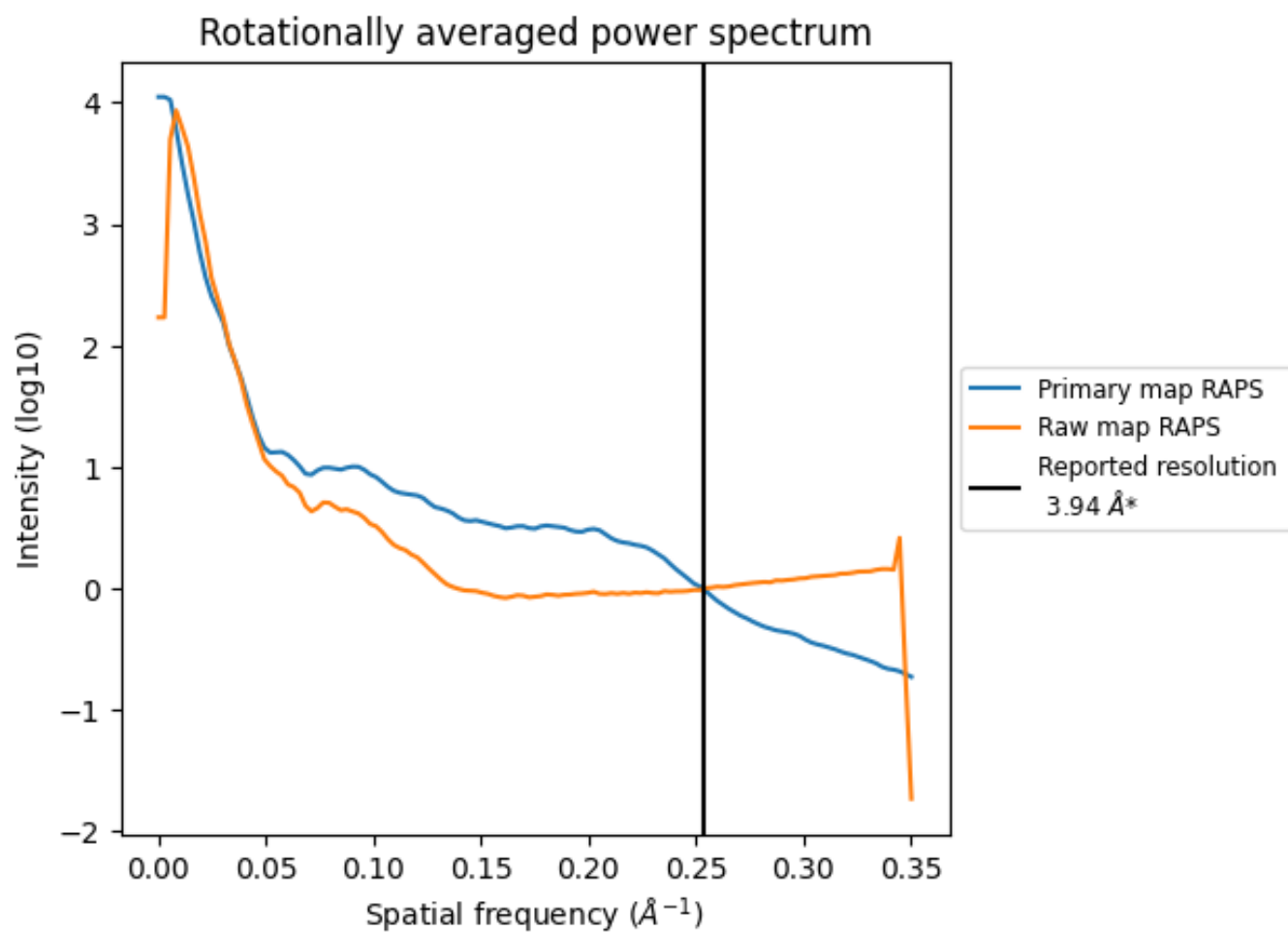
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 83 nm³; this corresponds to an approximate mass of 75 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

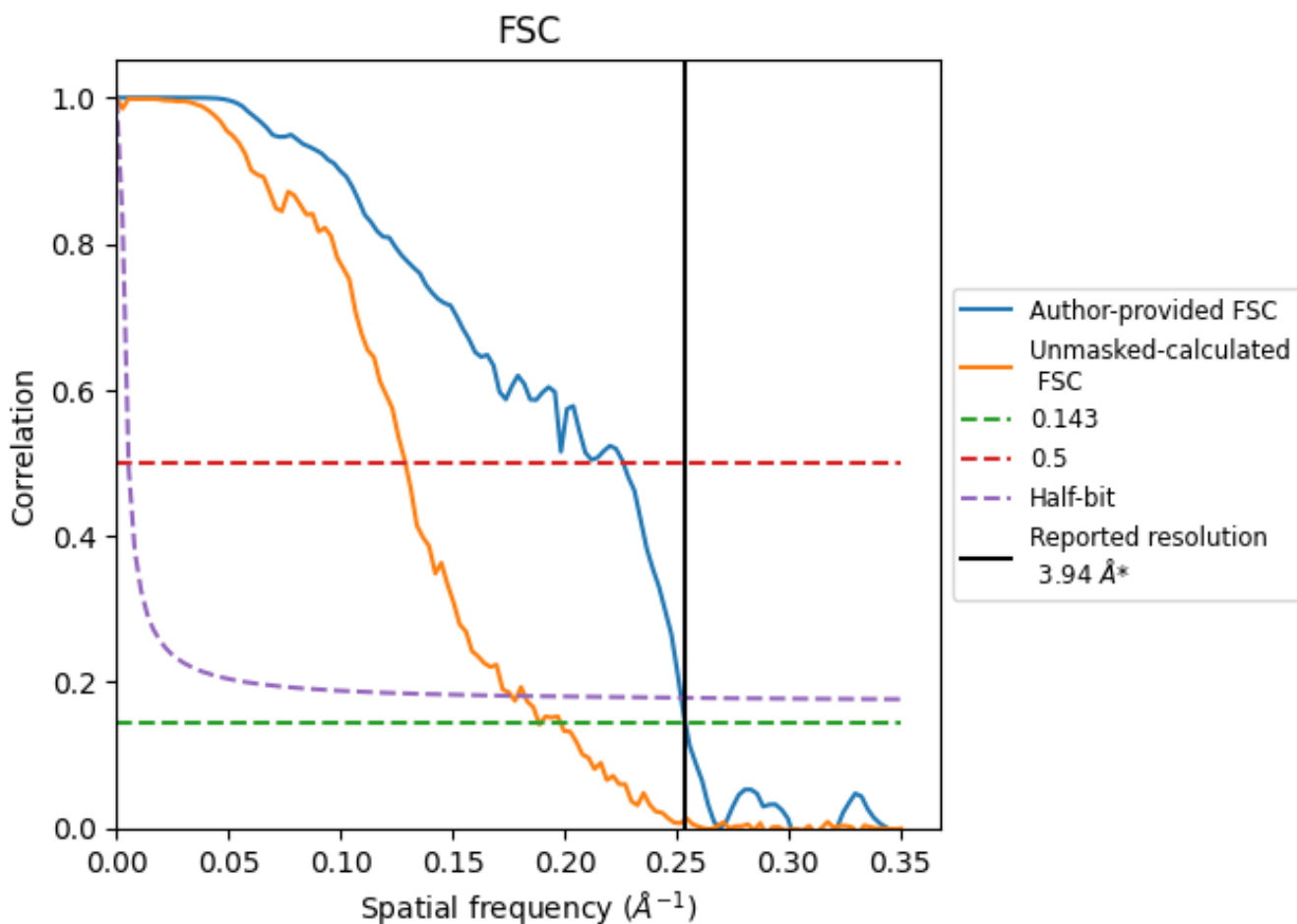


*Reported resolution corresponds to spatial frequency of 0.254 \AA^{-1}

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 0.254 Å⁻¹

8.2 Resolution estimates [i](#)

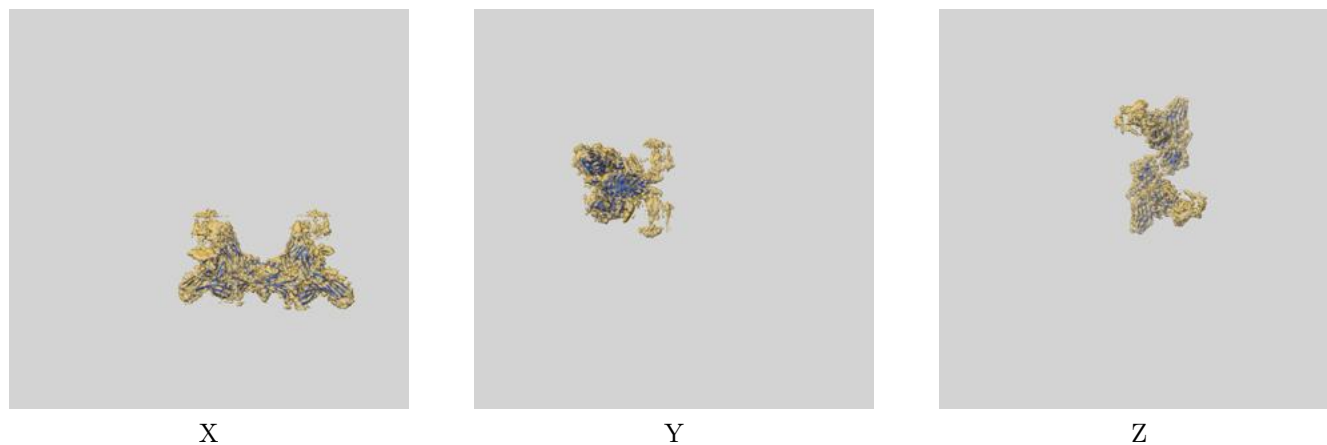
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.94	-	-
Author-provided FSC curve	3.94	4.42	3.97
Unmasked-calculated*	5.30	7.75	5.67

*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 5.30 differs from the reported value 3.94 by more than 10 %

9 Map-model fit [i](#)

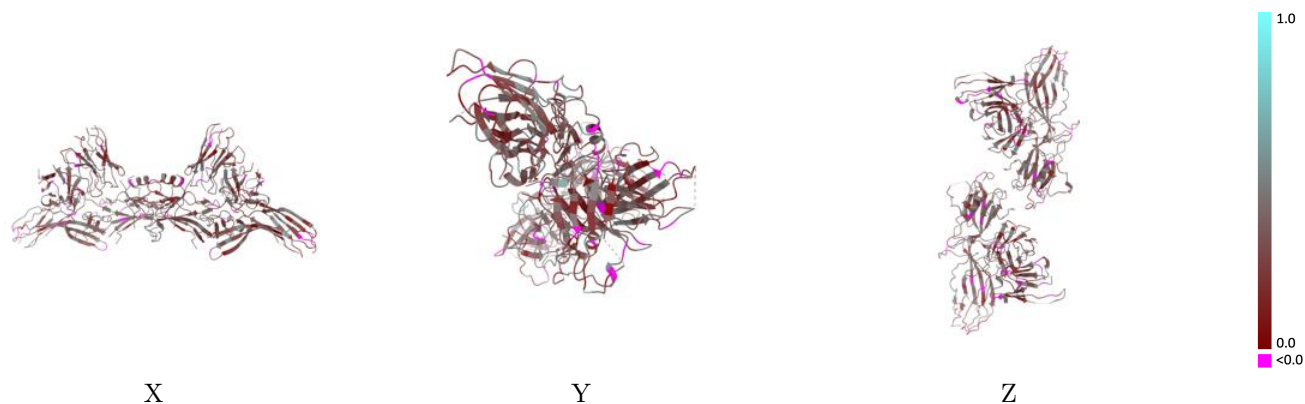
This section contains information regarding the fit between EMDB map EMD-36124 and PDB model 8JA3. Per-residue inclusion information can be found in section [3](#) on page [4](#).

9.1 Map-model overlay [i](#)



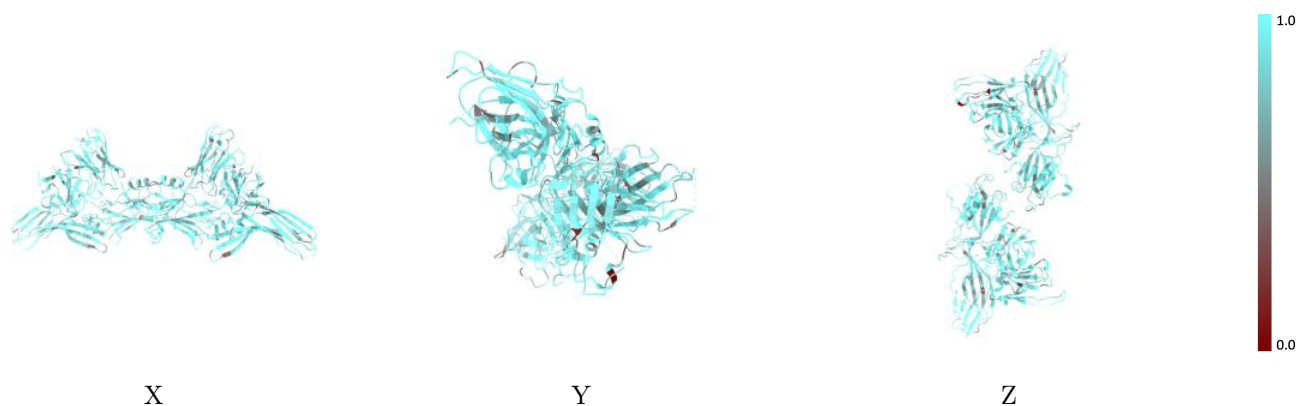
The images above show the 3D surface view of the map at the recommended contour level 0.044 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



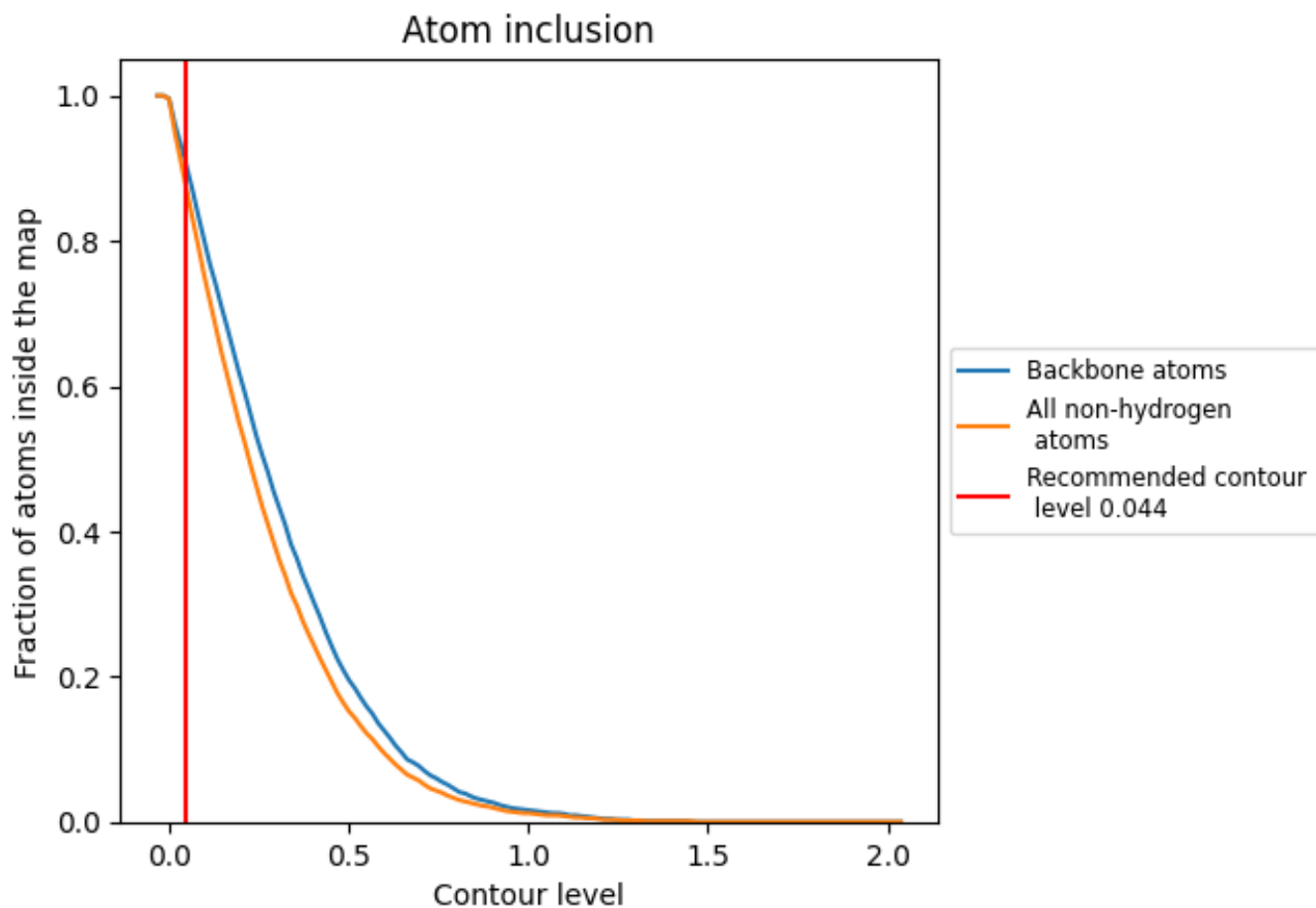
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.044).



















9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 88% 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 (0.044) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8780	 0.3270
A	 0.8870	 0.3310
B	 0.8830	 0.3260
H	 0.8650	 0.3210
I	 0.8540	 0.3170
L	 0.8740	 0.3210
M	 0.8620	 0.3170
U	 0.9140	 0.4000
V	 0.9630	 0.4320

