

wwPDB EM Validation Summary Report (i)

Jan 9, 2024 – 04:35 PM JST

PDB ID : 8J8Z EMDB ID : EMD-36082 Title Structure of beta-arrestin1 in complex with D6Rpp : Authors Maharana, J.; Sarma, P.; Yadav, M.K.; Chami, M.; Banerjee, R.; Shukla, A.K. : Deposited on 2023-05-02 : 3.40 Å(reported) Resolution : Based on initial model 8G08 :

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:

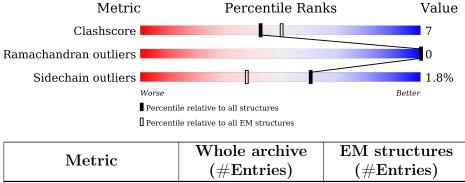
EMDB validation analysis	:	0.0.1.dev70
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al.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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.40 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	$(\# {\rm Entries})$	$(\# {\rm Entries})$
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	А	418	•		73%	8%	19%		
1	В	418	•		9%	19%			
2	Н	237		38%	11%	51%			
2	Ι	237		39%	9% •	51%			
3	L	215		43%	7%	50%			
3	М	215		39%	11%	50%			
4	U	18	11%	22%	6%	61%			
4	V	18	17%	17%	6%	61%			



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8722 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.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	А	340	Total 2601	C 1671		0 480		0	0
1	В	340	Total 2607	-		-	S 10	0	0

• Molecule 1 is a protein called Beta-arrestin-1.

• Molecule 2 is a protein called Fab30 Heavy Chain.

Mol	Chain	Residues	Atoms				AltConf	Trace	
2	Н	117	Total 898					0	0
2	Ι	117	Total 895	C 572				0	0

• Molecule 3 is a protein called Fab30 Light Chain.

Mol	Chain	Residues	Atoms				AltConf	Trace	
2	т	107	Total	С	Ν	0	S	0	0
0		107	804	508	132	161	3	0	U
9	М	107	Total	С	Ν	0	S	0	0
3	1/1	107	795	504	130	158	3	0	0

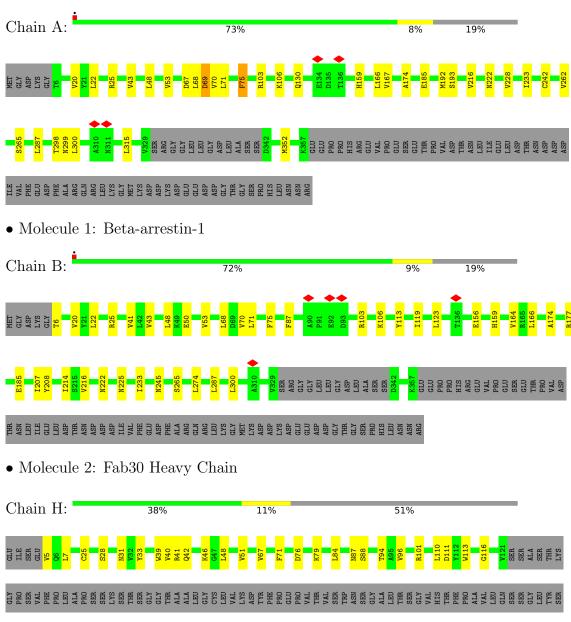
• Molecule 4 is a protein called Atypical chemokine receptor 2.

Mol	Chain	Residues	Atoms				AltConf	Trace		
4	V	7	Total	С	Ν	Ο	Р	S	0	0
4		1	61	29	7	21	3	1	0	0
4	II	7	Total	С	Ν	Ο	Р	S	0	0
4	U	1	61	29	7	21	3	1	0	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: Beta-arrestin-1



• Molecule 2: Fab30 Heavy Chain

Chain I	3	9%	9% •	51%
GLU ILE SER GLU V5 06	L7 V8 E9 E10 C11 L21 L21 L21 L23 L23	228 W39 W40 R41 K61 K68	NB0 184 184 189 189 189 189 193 193	D111 0115 0116 0116 0117 117 117 117 117 117 117 117 117 11
ALA PRO SER SER LYS SER	THR SER GLY GLY GLY ALA ALA ALA CVS CVS	LEU VAL LYS ASP ASP PHE PHE PRO GLU VAL	THR VAL SER TRP ASN ASN ALA ALA ALA ALA ALA ALA ALA VAL CLY VAL	PHE PRO ALA ALA ALA ALA CLN CLN CLN CLN CLN CLN CLN CLN CLN CLN
PRO SER SER LEU GLY	THR GLN THR TYR ILE CYS ASN VAL ASN	LYS PRO SER ASN THR LYS VAL ASP LYS LYS	VAL GLU PRO LYS SER CYS CYS CYS ASP LYS THR HIS HIS HIS	HIS HIS HIS HIS
• Molec	ule 3: Fab30	Light Chain		
Chain L	<i>i</i> :	43%	7%	50%
81 T6 L12 S13	Y37 Y37 Q39 K43 K43 L47 L47 L48	R62 D71 E82 B83 D83 091	V97 LTVS LTVS LTVS LTVS ARG ARG ARA ARA ARA ARA ARA ARA ARA ARA	PHE PHE PRO SER SER ASP SER CLN LLY CLN LLY SER SER ALA VAL VAL VAL VAL VAL ASN VAL
ASN PHE TYR PRO ARG GLU	ALA LYS VAL VAL GLN TRP CLN TRP LYS VAL ASN	LEU GLN SER ASN GLY GLU SER VAL	THR GLU GLN ASP SER LYS ASP ASP SER THR THR TYR SER SER SER	THR LEU THR THR SER SER LEU SER LYS CUU LYS CUU LYS CUU TYR VAL YAL VAL THR
HIS GLN GLY LEU SER SER	PRO VAL THR LYS SER PHE ASN ASN GLY	CAS		
• Molec	ule 3: Fab30	Light Chain		
Chain N	И :	39%	11%	50%
S1 D2 Q4 M5	L112 S13 Y37 Y40 F41 L47 L47	149 851 851 852 756 756 756 756 756	L74 E82 B83 B83 C89 Q90 Q91 Q91 Q91 C89 Q100	1103 LIVS LIVS LIVS LIVS ARR ARA ALA ALA ALA ARA PRO PRO PRO PRO SER SER SER
GLN LEU LYS SER GLY THR	ALA SER VAL VAL CYS CYS LEU LEU ASN ASN	TYR PRO ARG GLU ALA LYS CLN TRP LYS	VAL ASP ASP ALA ALA CLU SER GLN SER CLU SER VAL	THR CLU GLU GLU GLU GLU GLU LYS SER SER THR THR THR THR THR THR THR THR THR TH
ALA ASP TYR GLU LYS HIS	LYS VAL TYR ALA ALA CYS GLU VAL THR HIS GLN	GLY GLY LEU SER SER PRO VAL THR THR LYS SER PHE	ASN ARG GLY GLU CYS	
• Molec	ule 4: Atypica	al chemokine r	eceptor 2	
Chain V	/: 17%	17% 6%		51%
GLY TPO ALA GLN ALA SEP	LEU SEP SEP SEP C347 C347 C347 E349 S351 I 355 C355 L355	ALA		
• Molec	ule 4: Atypica	al chemokine r	eceptor 2	
Chain U	J: 11%	22% 6%	6	1%







4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	369871	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 $(e^-/\text{\AA}^2)$	53	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	2.437	Depositor
Minimum map value	-1.250	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.025	Depositor
Recommended contour level	0.15	Depositor
Map size (Å)	421.4304, 421.4304, 421.4304	wwPDB
Map dimensions	288, 288, 288	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.4633, 1.4633, 1.4633	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: 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		
IVIOI	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.29	0/2657	0.51	0/3618	
1	В	0.28	0/2664	0.50	0/3627	
2	Н	0.34	0/922	0.55	0/1255	
2	Ι	0.34	0/919	0.56	0/1251	
3	L	0.33	0/822	0.51	0/1119	
3	М	0.35	0/813	0.57	0/1108	
4	U	0.57	0/28	0.87	0/33	
4	V	0.50	0/28	0.63	0/33	
All	All	0.31	0/8853	0.52	0/12044	

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (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	А	2601	0	2598	32	0
1	В	2607	0	2618	29	0
2	Н	898	0	845	21	0
2	Ι	895	0	841	16	0
3	L	804	0	776	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	М	795	0	761	17	0
4	U	61	0	42	4	0
4	V	61	0	42	8	0
All	All	8722	0	8523	123	0

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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 7.

The worst 5 of 123 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:50:GLU:N	1:B:50:GLU:OE1	2.12	0.82
1:B:103:ARG:HD2	4:U:353:LEU:HD22	1.65	0.77
1:B:216:VAL:HG21	1:B:233:ILE:HD11	1.69	0.72
1:B:48:LEU:HD11	1:B:53:VAL:HG23	1.74	0.69
3:L:82:GLU:OE1	3:L:82:GLU:N	2.24	0.69

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	А	336/418~(80%)	312 (93%)	24 (7%)	0	100 100
1	В	336/418~(80%)	313~(93%)	23 (7%)	0	100 100
2	Н	115/237~(48%)	108 (94%)	7 (6%)	0	100 100
2	Ι	115/237~(48%)	109~(95%)	6 (5%)	0	100 100
3	L	105/215~(49%)	96 (91%)	9 (9%)	0	100 100
3	М	105/215~(49%)	99~(94%)	6 (6%)	0	100 100
4	U	2/18~(11%)	1 (50%)	1 (50%)	0	100 100

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	3	Analysed	Favoured	Allowed	Outliers	Percentiles
4	V	2/18~(11%)	1 (50%)	1 (50%)	0	100 100
All	All	1116/1776~(63%)	1039~(93%)	77 (7%)	0	100 100

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There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	А	281/372~(76%)	277~(99%)	4 (1%)	67 83
1	В	282/372~(76%)	279~(99%)	3~(1%)	73 86
2	Н	92/200~(46%)	91~(99%)	1 (1%)	73 86
2	Ι	91/200~(46%)	86~(94%)	5~(6%)	21 51
3	L	89/190~(47%)	88~(99%)	1 (1%)	73 86
3	М	86/190~(45%)	85~(99%)	1 (1%)	71 85
4	U	4/6~(67%)	3~(75%)	1 (25%)	0 2
4	V	4/6~(67%)	3~(75%)	1 (25%)	0 2
All	All	929/1536~(60%)	912~(98%)	17~(2%)	61 79

5 of 17 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
2	Ι	107	TYR
4	U	347	CYS
1	В	75	PHE
1	В	177	ARG
1	В	225	ASN

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such side chains are listed below:



Mol	Chain	Res	Type
1	А	159	HIS
2	Н	80	ASN
2	Ι	42	GLN
3	М	39	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)

6 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	Trune	Chain	Res	Link	B	ond leng	gths	В	ond ang	gles
NIOI	Type	Unam	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
4	SEP	V	350	4	8,9,10	1.52	1 (12%)	8,12,14	1.25	1 (12%)
4	SEP	V	348	4	8,9,10	0.56	0	8,12,14	0.66	0
4	SEP	V	351	4	8,9,10	0.63	0	8,12,14	0.64	0
4	SEP	U	350	4	8,9,10	1.51	1 (12%)	8,12,14	1.29	1 (12%)
4	SEP	U	348	4	8,9,10	1.53	1 (12%)	8,12,14	1.56	2 (25%)
4	SEP	U	351	4	8,9,10	0.62	0	8,12,14	0.65	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SEP	V	350	4	-	1/5/8/10	-
4	SEP	V	348	4	-	1/5/8/10	-
4	SEP	V	351	4	-	4/5/8/10	-
4	SEP	U	350	4	-	0/5/8/10	-
4	SEP	U	348	4	-	0/5/8/10	-
4	SEP	U	351	4	-	4/5/8/10	-



Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
4	U	348	SEP	P-01P	3.33	1.61	1.50
4	U	350	SEP	P-O1P	3.31	1.61	1.50
4	V	350	SEP	P-O1P	3.31	1.61	1.50

All (3) bond length outliers are listed below:

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
4	U	348	SEP	OG-CB-CA	2.82	110.89	108.14
4	V	350	SEP	P-OG-CB	-2.74	110.74	118.30
4	U	348	SEP	P-OG-CB	-2.74	110.76	118.30
4	U	350	SEP	P-OG-CB	-2.52	111.36	118.30

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	V	348	SEP	N-CA-CB-OG
4	V	351	SEP	CB-OG-P-O2P
4	V	351	SEP	CB-OG-P-O3P
4	U	351	SEP	N-CA-CB-OG
4	U	351	SEP	CB-OG-P-O1P

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	V	348	SEP	1	0
4	U	348	SEP	1	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.



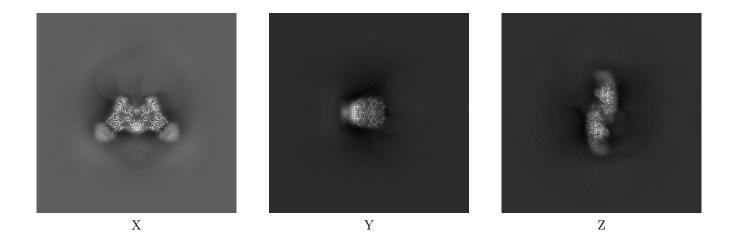
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-36082. These allow visual inspection of the internal detail of the map and identification of artifacts.

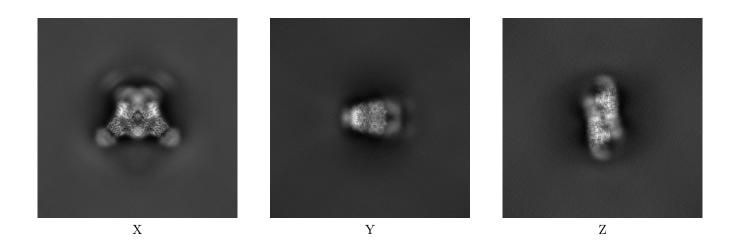
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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.

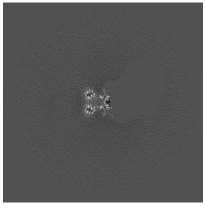


6.2 Central slices (i)

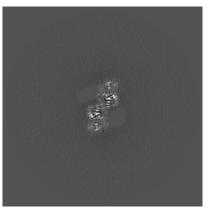
6.2.1 Primary map



X Index: 144

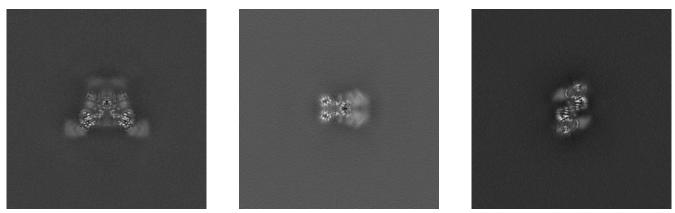


Y Index: 144



Z Index: 144

6.2.2 Raw map



X Index: 144

Y Index: 144

Z Index: 144

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

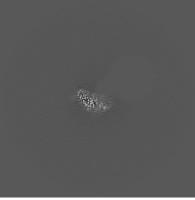


6.3 Largest variance slices (i)

6.3.1 Primary map



X Index: 136

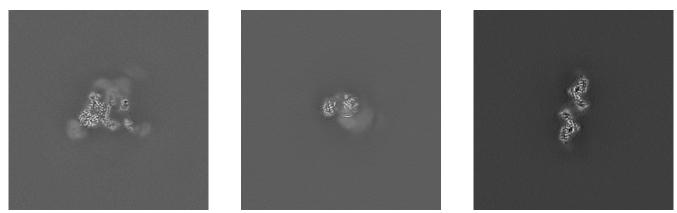


Y Index: 114



Z Index: 130

6.3.2 Raw map



X Index: 136

Y Index: 168

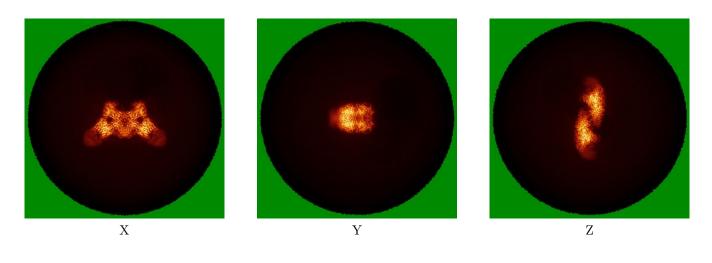


The images above show the largest variance slices of the map in three orthogonal directions.

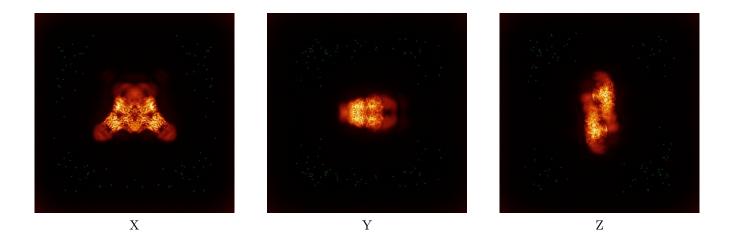


6.4 Orthogonal standard-deviation projections (False-color) (i)

6.4.1 Primary map



6.4.2 Raw map

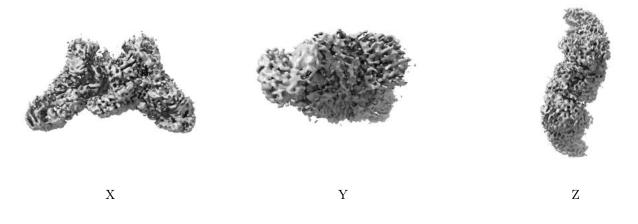


The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.



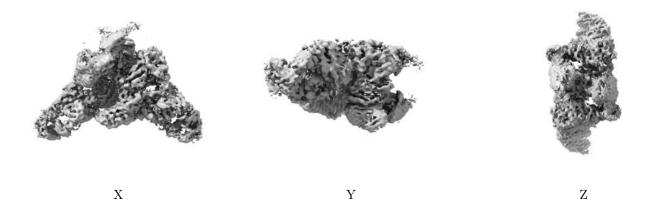
6.5 Orthogonal surface views (i)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.15. 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)

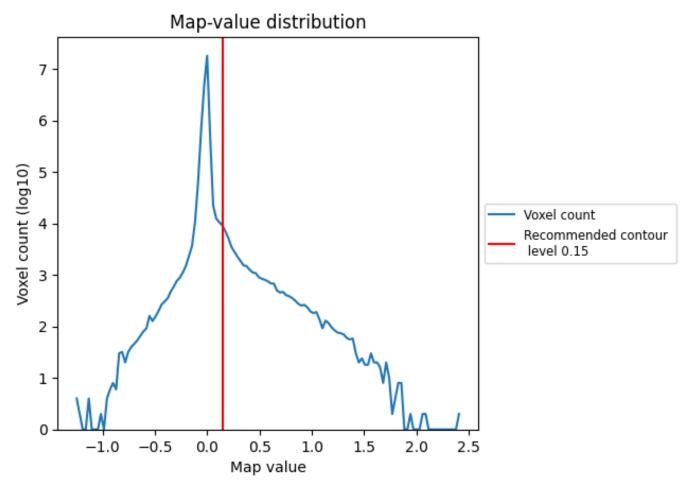
This section was not generated. No masks/segmentation were deposited.



7 Map analysis (i)

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

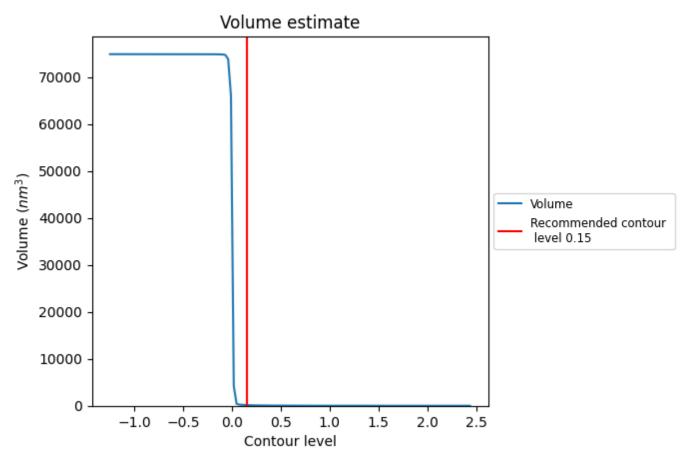
7.1 Map-value distribution (i)



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)

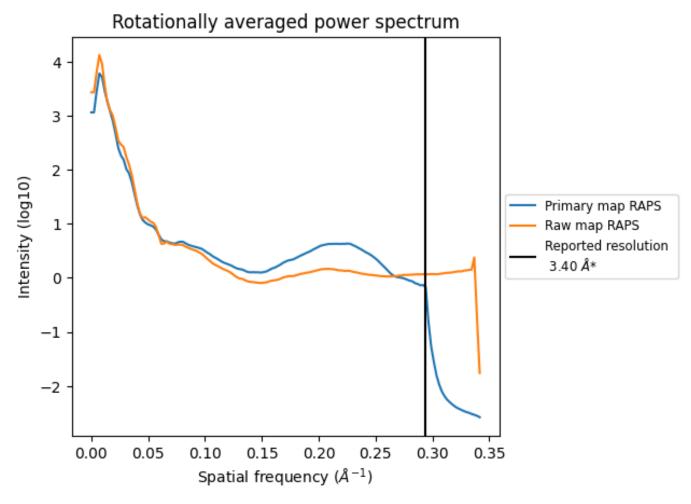


The volume at the recommended contour level is 147 $\rm nm^3;$ this corresponds to an approximate mass of 132 kDa.

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)



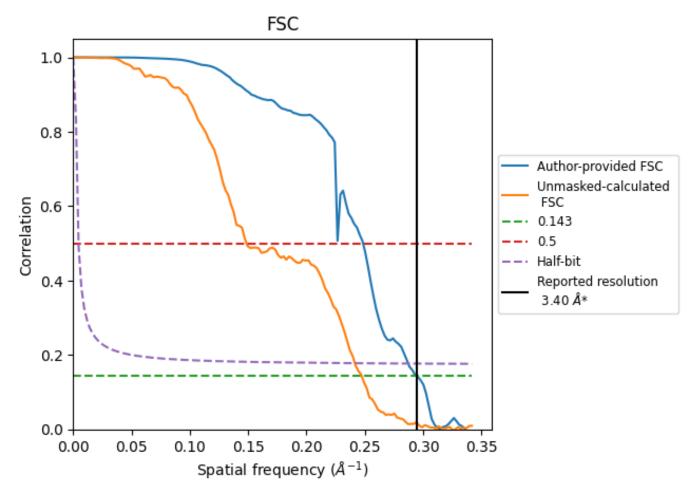
*Reported resolution corresponds to spatial frequency of 0.294 $\mathrm{\AA^{-1}}$



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 0.294 $\mathrm{\AA^{-1}}$



8.2 Resolution estimates (i)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
Resolution estimate (A)	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	3.40	4.03	3.48
Unmasked-calculated*	4.05	6.73	4.14

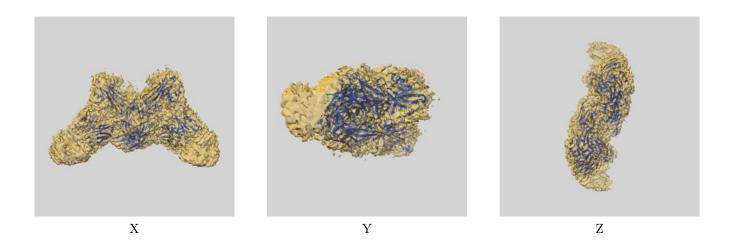
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.05 differs from the reported value 3.4 by more than 10 %



9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-36082 and PDB model 8J8Z. Per-residue inclusion information can be found in section 3 on page 4.

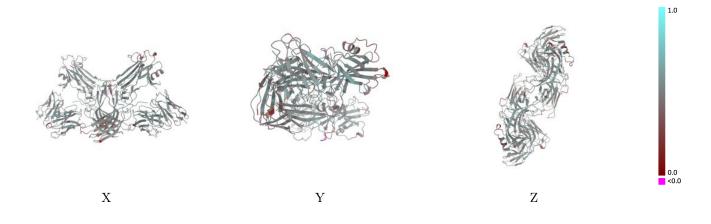
9.1 Map-model overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.15 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.

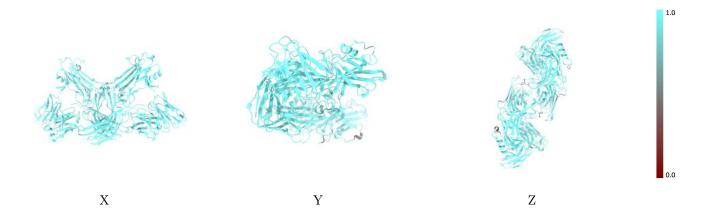


9.2 Q-score mapped to coordinate model (i)



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.

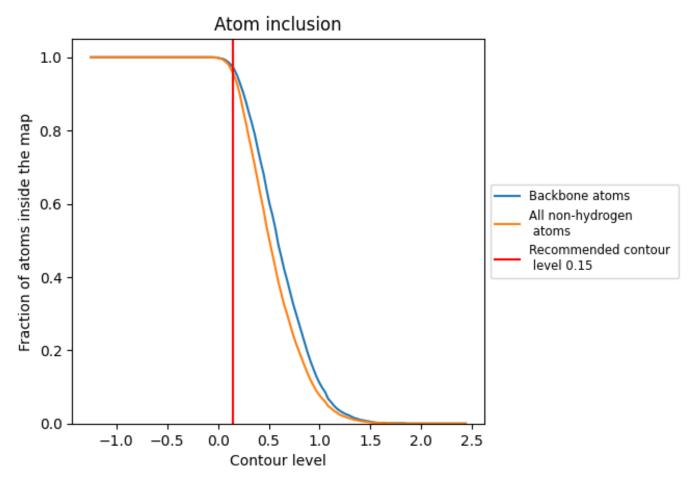
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.15).



9.4 Atom inclusion (i)



At the recommended contour level, 97% of all backbone atoms, 96% 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 (0.15) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9560	0.4890
А	0.9420	0.4760
В	0.9440	0.4790
Н	0.9700	0.5070
Ι	0.9750	0.5110
L	0.9800	0.5020
М	0.9760	0.4970
U	0.9840	0.5290
V	1.0000	0.5270

