



Full wwPDB EM Validation Report ⓘ

Mar 24, 2026 – 02:27 PM UTC

PDB ID : 8EXQ / pdb_00008exq
EMDB ID : EMD-28659
Title : Cryo-EM structure of S. aureus BlaR1 with C1 symmetry
Authors : Worrall, L.J.; Alexander, J.A.N.; Vuckovic, M.; Strynadka, N.C.J.
Deposited on : 2022-10-25
Resolution : 4.90 Å(reported)

This is a Full wwPDB EM Validation 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.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

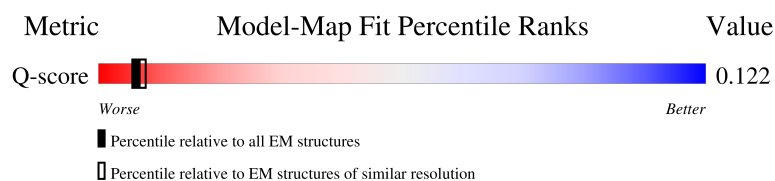
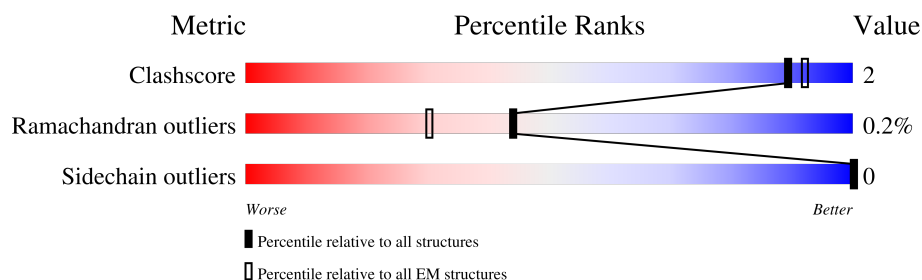
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.90 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	1274 (4.40 - 5.40)

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	602	28% 84% 7% 8%
1	B	602	23% 82% 10% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9304 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-lactam sensor/signal transducer BlaR1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	554	Total	C	N	O	S	0	0
			4651	3056	767	809	19		
1	A	554	Total	C	N	O	S	0	0
			4651	3056	767	809	19		

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	586	GLU	-	expression tag	UNP Q00419
B	587	LEU	-	expression tag	UNP Q00419
B	588	ALA	-	expression tag	UNP Q00419
B	589	LEU	-	expression tag	UNP Q00419
B	590	VAL	-	expression tag	UNP Q00419
B	591	PRO	-	expression tag	UNP Q00419
B	592	ARG	-	expression tag	UNP Q00419
B	593	GLY	-	expression tag	UNP Q00419
B	594	SER	-	expression tag	UNP Q00419
B	595	SER	-	expression tag	UNP Q00419
B	596	ALA	-	expression tag	UNP Q00419
B	597	HIS	-	expression tag	UNP Q00419
B	598	HIS	-	expression tag	UNP Q00419
B	599	HIS	-	expression tag	UNP Q00419
B	600	HIS	-	expression tag	UNP Q00419
B	601	HIS	-	expression tag	UNP Q00419
B	602	HIS	-	expression tag	UNP Q00419
A	586	GLU	-	expression tag	UNP Q00419
A	587	LEU	-	expression tag	UNP Q00419
A	588	ALA	-	expression tag	UNP Q00419
A	589	LEU	-	expression tag	UNP Q00419
A	590	VAL	-	expression tag	UNP Q00419
A	591	PRO	-	expression tag	UNP Q00419
A	592	ARG	-	expression tag	UNP Q00419
A	593	GLY	-	expression tag	UNP Q00419
A	594	SER	-	expression tag	UNP Q00419

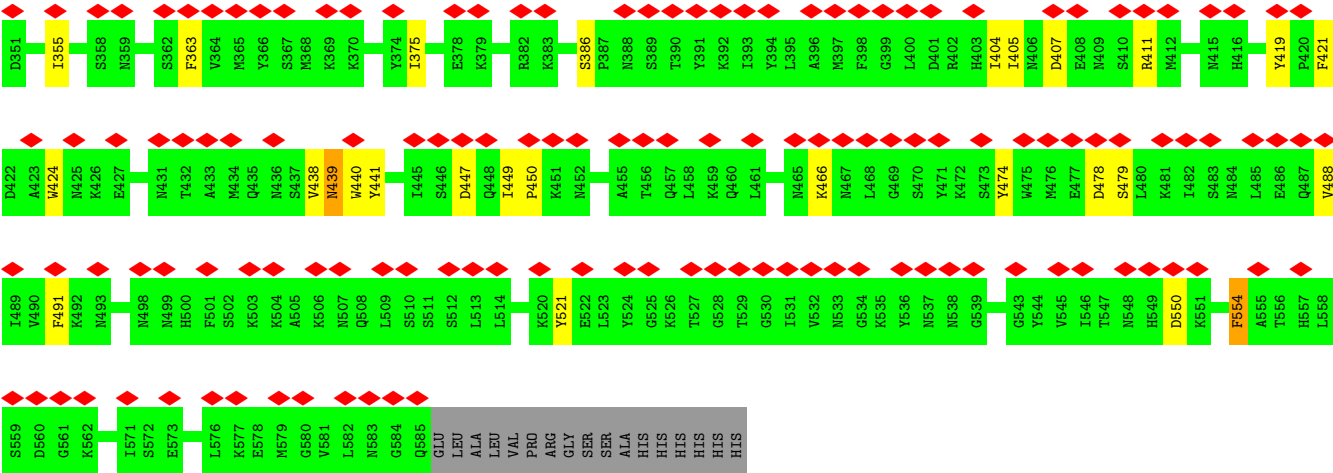
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Chain	Residue	Modelled	Actual	Comment	Reference
A	595	SER	-	expression tag	UNP Q00419
A	596	ALA	-	expression tag	UNP Q00419
A	597	HIS	-	expression tag	UNP Q00419
A	598	HIS	-	expression tag	UNP Q00419
A	599	HIS	-	expression tag	UNP Q00419
A	600	HIS	-	expression tag	UNP Q00419
A	601	HIS	-	expression tag	UNP Q00419
A	602	HIS	-	expression tag	UNP Q00419

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
2	B	1	Total 1	Zn 1	0
2	A	1	Total 1	Zn 1	0



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	23797	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	1.035	Depositor
Minimum map value	-0.607	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.045	Depositor
Recommended contour level	0.17	Depositor
Map size (Å)	231.0016, 231.0016, 231.0016	wwPDB
Map dimensions	128, 128, 128	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.8047, 1.8047, 1.8047	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

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	1.09	1/4766 (0.0%)	1.23	36/6418 (0.6%)
1	B	1.10	4/4766 (0.1%)	1.26	43/6418 (0.7%)
All	All	1.09	5/9532 (0.1%)	1.25	79/12836 (0.6%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	201	HIS	CB-CG	-7.17	1.40	1.50
1	B	153	HIS	CB-CG	-5.68	1.42	1.50
1	B	552	TYR	CB-CG	-5.25	1.40	1.51
1	A	90	PHE	CB-CG	-5.08	1.39	1.50
1	B	494	MET	CG-SD	-5.02	1.68	1.80

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	PHE	CA-CB-CG	-7.97	105.83	113.80
1	A	550	ASP	CA-CB-CG	7.97	120.57	112.60
1	B	386	SER	CA-C-N	7.83	127.47	119.56
1	B	386	SER	C-N-CA	7.83	127.47	119.56
1	B	419	TYR	CA-C-N	7.82	127.53	119.56
1	B	419	TYR	C-N-CA	7.82	127.53	119.56
1	A	419	TYR	CA-C-N	7.72	127.44	119.56
1	A	419	TYR	C-N-CA	7.72	127.44	119.56
1	B	479	SER	N-CA-C	7.51	120.41	111.33
1	A	386	SER	CA-C-N	7.42	127.12	119.56
1	A	386	SER	C-N-CA	7.42	127.12	119.56
1	B	550	ASP	CA-CB-CG	7.37	119.97	112.60
1	A	338	ASN	CA-CB-CG	-7.32	105.28	112.60
1	A	405	ILE	N-CA-C	-7.14	98.20	108.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	272	VAL	CA-C-N	7.10	127.06	120.03
1	B	272	VAL	C-N-CA	7.10	127.06	120.03
1	B	491	PHE	CA-CB-CG	-7.08	106.72	113.80
1	A	479	SER	N-CA-C	7.04	119.85	111.33
1	B	341	LYS	CA-C-N	6.98	126.91	120.21
1	B	341	LYS	C-N-CA	6.98	126.91	120.21
1	A	236	LYS	N-CA-C	-6.78	103.35	111.69
1	A	271	ASN	N-CA-C	6.78	118.67	111.28
1	B	227	ASN	CA-C-N	6.63	126.33	119.56
1	B	227	ASN	C-N-CA	6.63	126.33	119.56
1	B	375	ILE	N-CA-C	6.55	117.31	107.75
1	A	227	ASN	CA-C-N	6.54	127.06	119.47
1	A	227	ASN	C-N-CA	6.54	127.06	119.47
1	B	5	LEU	N-CA-C	-6.38	103.85	111.69
1	B	442	PHE	CA-CB-CG	-6.35	107.45	113.80
1	A	272	VAL	CA-C-N	6.31	126.28	120.03
1	A	272	VAL	C-N-CA	6.31	126.28	120.03
1	B	313	PHE	CA-CB-CG	-6.29	107.50	113.80
1	A	253	ILE	N-CA-C	6.29	116.91	108.11
1	B	404	ILE	N-CA-C	-6.28	103.03	111.44
1	B	12	PHE	CA-CB-CG	6.11	119.91	113.80
1	B	478	ASP	CA-C-N	6.10	128.74	120.38
1	B	478	ASP	C-N-CA	6.10	128.74	120.38
1	A	447	ASP	CA-CB-CG	6.08	118.68	112.60
1	B	174	TRP	N-CA-C	6.08	118.44	111.02
1	B	187	PHE	N-CA-C	-6.07	104.95	112.90
1	B	160	VAL	N-CA-C	6.05	116.58	108.11
1	A	439	ASN	CA-CB-CG	5.88	118.48	112.60
1	B	417	LYS	N-CA-C	-5.85	101.86	110.52
1	A	363	PHE	N-CA-C	5.84	118.02	108.55
1	A	478	ASP	CA-C-N	5.83	128.37	120.38
1	A	478	ASP	C-N-CA	5.83	128.37	120.38
1	B	150	LEU	N-CA-C	5.75	120.42	113.17
1	A	272	VAL	N-CA-C	5.71	115.30	108.45
1	B	489	ILE	CA-C-N	5.70	127.85	120.56
1	B	489	ILE	C-N-CA	5.70	127.85	120.56
1	B	447	ASP	CA-CB-CG	5.69	118.29	112.60
1	A	248	PHE	CA-CB-CG	5.69	119.49	113.80
1	A	411	ARG	N-CA-C	5.65	118.18	110.55
1	A	554	PHE	N-CA-C	5.62	117.39	108.79
1	A	404	ILE	N-CA-C	-5.58	103.97	111.44
1	B	328	PHE	CA-CB-CG	-5.57	108.23	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	ILE	N-CA-C	5.56	116.11	107.28
1	B	411	ARG	N-CA-C	5.49	119.05	110.32
1	A	355	ILE	N-CA-C	-5.49	106.35	111.45
1	B	93	ASP	N-CA-C	5.48	117.68	108.96
1	A	491	PHE	CA-CB-CG	-5.47	108.33	113.80
1	A	341	LYS	CA-C-N	5.46	125.93	120.14
1	A	341	LYS	C-N-CA	5.46	125.93	120.14
1	B	473	SER	CA-C-N	5.36	129.26	120.63
1	B	473	SER	C-N-CA	5.36	129.26	120.63
1	B	529	THR	N-CA-C	5.31	118.10	109.24
1	A	466	LYS	N-CA-CB	-5.28	104.42	112.28
1	B	540	TRP	CA-CB-CG	5.27	123.61	113.60
1	B	232	ILE	N-CA-C	-5.21	106.61	111.45
1	B	501	PHE	CA-CB-CG	5.20	119.00	113.80
1	B	363	PHE	N-CA-C	5.17	116.92	108.55
1	A	333	SER	CA-C-N	5.11	129.00	120.62
1	A	333	SER	C-N-CA	5.11	129.00	120.62
1	B	188	LYS	N-CA-C	-5.09	105.62	111.07
1	B	427	GLU	N-CA-C	-5.09	103.68	110.55
1	A	152	ASN	CA-CB-CG	-5.05	107.55	112.60
1	B	160	VAL	CA-C-N	-5.04	116.56	123.12
1	B	160	VAL	C-N-CA	-5.04	116.56	123.12
1	A	375	ILE	N-CA-CB	-5.02	105.57	111.90

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4651	0	4747	15	0
1	B	4651	0	4747	19	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
All	All	9304	0	9494	34	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (34) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:THR:HG22	1:B:282:HIS:CD2	2.12	0.84
1:B:211:THR:CG2	1:B:282:HIS:CD2	2.61	0.83
1:B:211:THR:HG23	1:B:282:HIS:HD2	1.52	0.73
1:B:211:THR:HG23	1:B:282:HIS:CD2	2.30	0.65
1:A:175:TYR:CG	1:A:175:TYR:O	2.56	0.57
1:B:362:SER:OG	1:B:484:ASN:ND2	2.39	0.55
1:B:336:ASP:OD1	1:B:481:LYS:NZ	2.41	0.54
1:B:211:THR:CG2	1:B:282:HIS:HD2	2.07	0.53
1:B:152:ASN:N	1:B:152:ASN:OD1	2.41	0.51
1:B:407:ASP:OD2	1:B:504:LYS:NZ	2.44	0.49
1:B:151:PHE:CD1	1:B:151:PHE:C	2.90	0.49
1:A:421:PHE:HB2	1:A:424:TRP:CD1	2.49	0.48
1:B:166:THR:OG1	1:B:167:ILE:N	2.48	0.47
1:B:426:LYS:O	1:B:428:GLN:NE2	2.48	0.47
1:A:152:ASN:OD1	1:A:152:ASN:N	2.39	0.47
1:A:158:ASN:N	1:A:158:ASN:OD1	2.47	0.47
1:B:527:THR:OG1	1:B:528:GLY:N	2.49	0.46
1:A:192:ASP:CG	1:A:193:LYS:N	2.75	0.45
1:A:407:ASP:OD1	1:A:407:ASP:N	2.48	0.45
1:A:274:PHE:CD1	1:A:274:PHE:C	2.91	0.45
1:B:274:PHE:CD1	1:B:274:PHE:C	2.94	0.45
1:A:438:VAL:HG12	1:A:441:TYR:H	1.81	0.45
1:A:439:ASN:O	1:A:440:TRP:C	2.62	0.42
1:B:175:TYR:O	1:B:175:TYR:CG	2.68	0.42
1:B:137:TYR:CD2	1:B:137:TYR:C	2.96	0.42
1:A:449:ILE:O	1:A:450:PRO:C	2.63	0.41
1:A:88:GLN:HB2	1:A:90:PHE:CE1	2.55	0.41
1:A:521:TYR:CD1	1:A:521:TYR:C	2.96	0.41
1:A:488:VAL:HG22	1:A:554:PHE:HA	2.01	0.41
1:A:424:TRP:CD1	1:A:438:VAL:HG22	2.56	0.41
1:B:122:LEU:O	1:B:123:LYS:C	2.63	0.41
1:A:147:ASP:O	1:A:151:PHE:N	2.54	0.41
1:B:438:VAL:HG12	1:B:441:TYR:H	1.86	0.40
1:B:174:TRP:O	1:B:176:GLY:N	2.53	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	550/602 (91%)	536 (98%)	13 (2%)	1 (0%)	43	77
1	B	550/602 (91%)	534 (97%)	15 (3%)	1 (0%)	43	77
All	All	1100/1204 (91%)	1070 (97%)	28 (2%)	2 (0%)	44	77

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	175	TYR
1	A	474	TYR

5.3.2 Protein sidechains

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	518/562 (92%)	518 (100%)	0	100	100
1	B	518/562 (92%)	518 (100%)	0	100	100
All	All	1036/1124 (92%)	1036 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	239	HIS
1	B	282	HIS

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Mol	Chain	Res	Type
1	B	416	HIS
1	B	431	ASN
1	B	484	ASN
1	B	508	GLN
1	A	285	ASN
1	A	325	GLN
1	A	327	GLN
1	A	403	HIS
1	A	448	GLN
1	A	538	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](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are 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 ⓘ

There are no chain breaks in this entry.

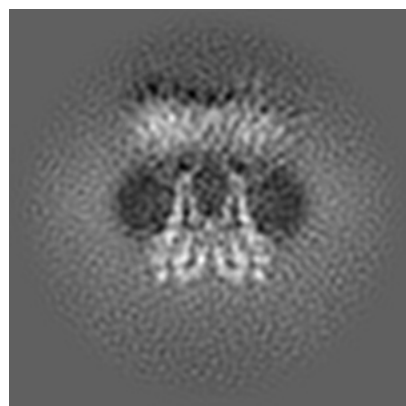
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-28659. These allow visual inspection of the internal detail of the map and identification of artifacts.

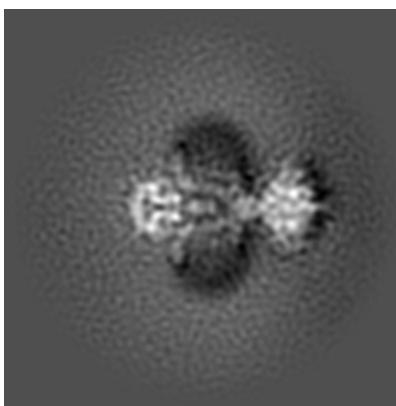
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

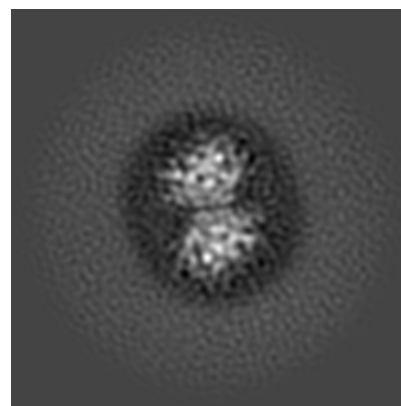
6.1.1 Primary map



X

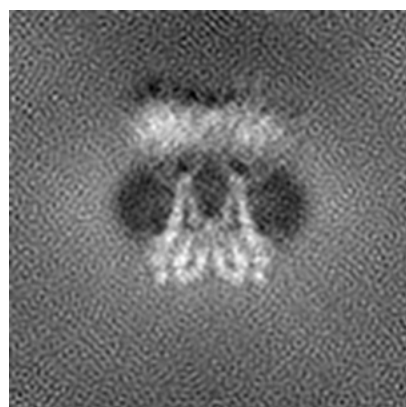


Y

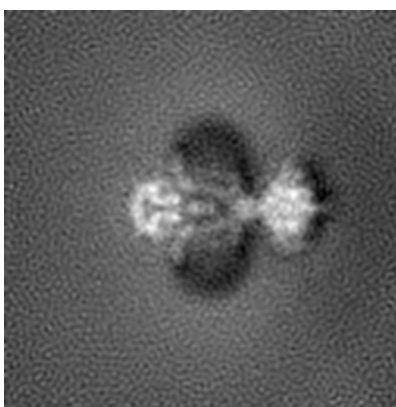


Z

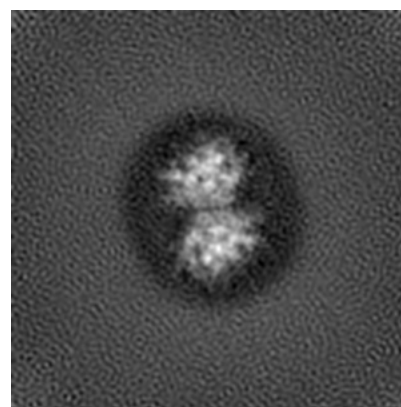
6.1.2 Raw map



X



Y

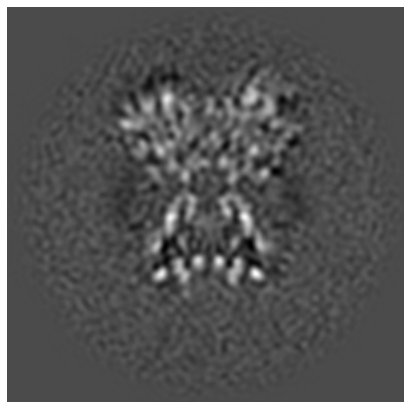


Z

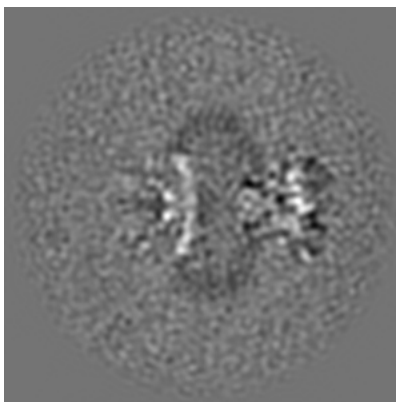
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

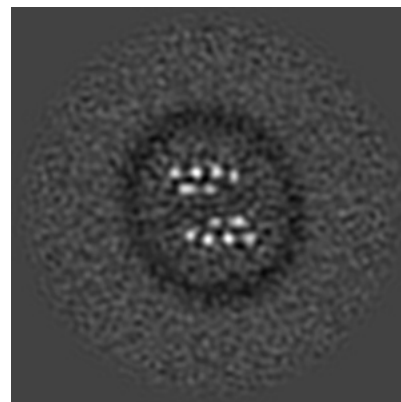
6.2.1 Primary map



X Index: 64

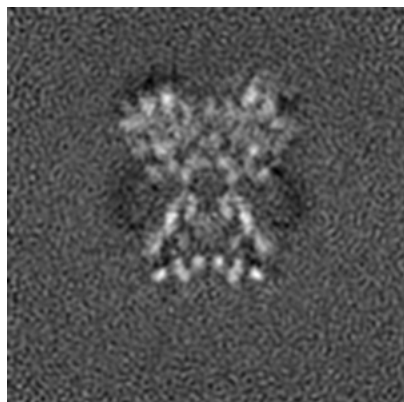


Y Index: 64

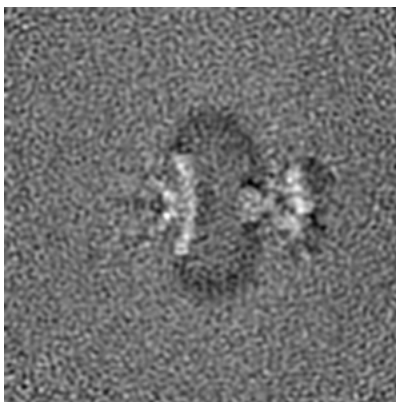


Z Index: 64

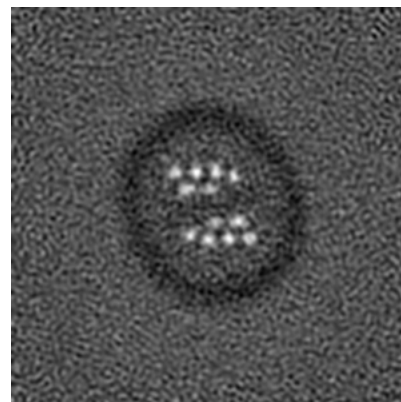
6.2.2 Raw map



X Index: 64



Y Index: 64

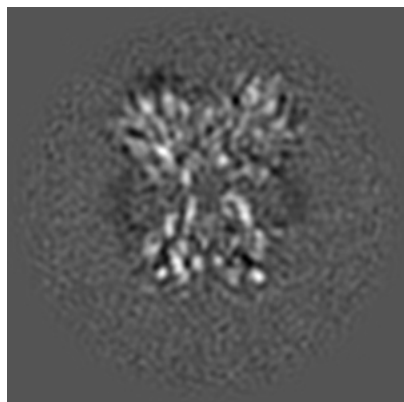


Z Index: 64

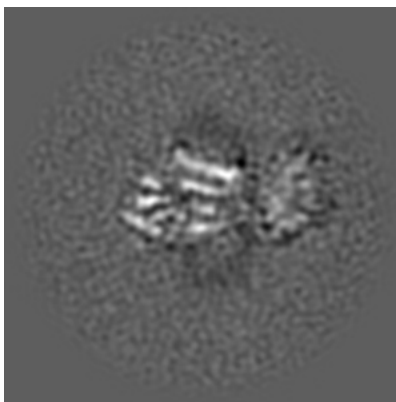
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

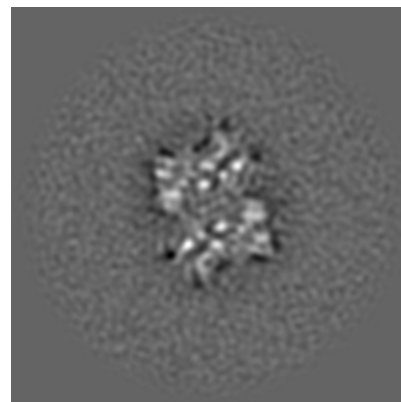
6.3.1 Primary map



X Index: 65

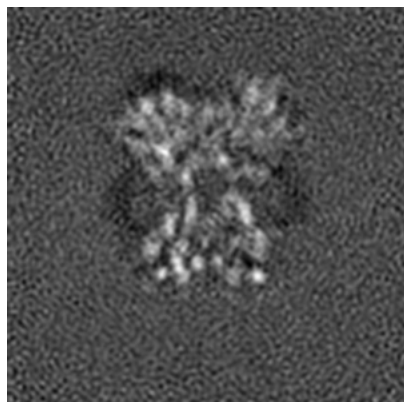


Y Index: 54

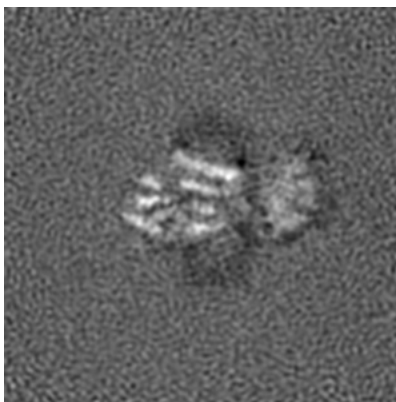


Z Index: 56

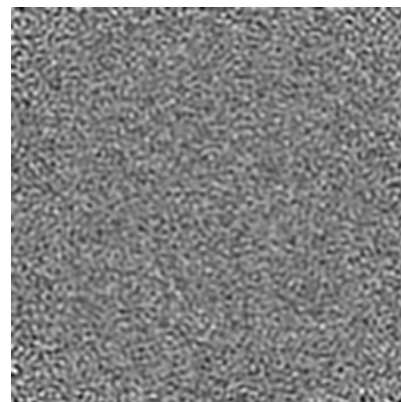
6.3.2 Raw map



X Index: 65



Y Index: 54

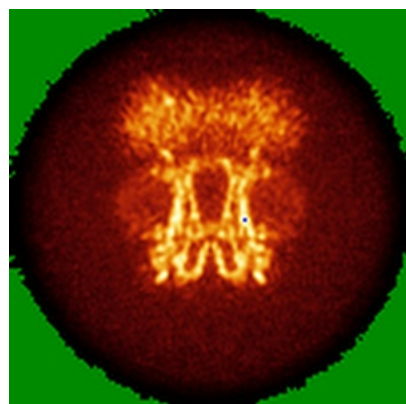


Z Index: 0

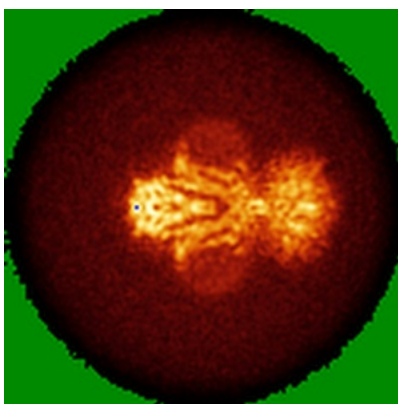
The images above show the largest variance slices of the map in three orthogonal directions.

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

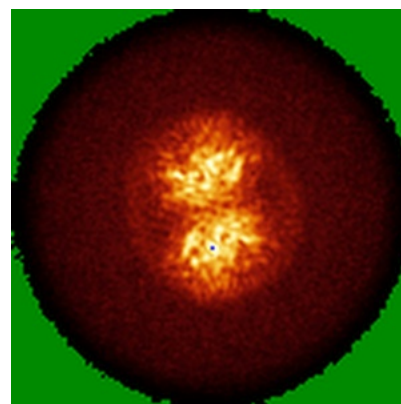
6.4.1 Primary map



X

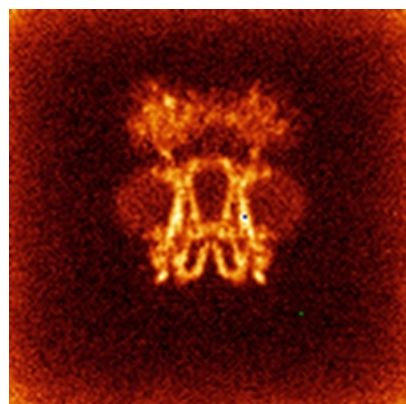


Y

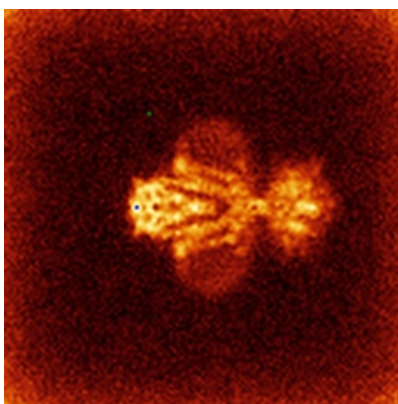


Z

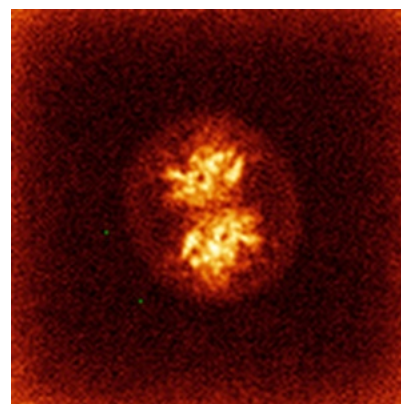
6.4.2 Raw map



X



Y

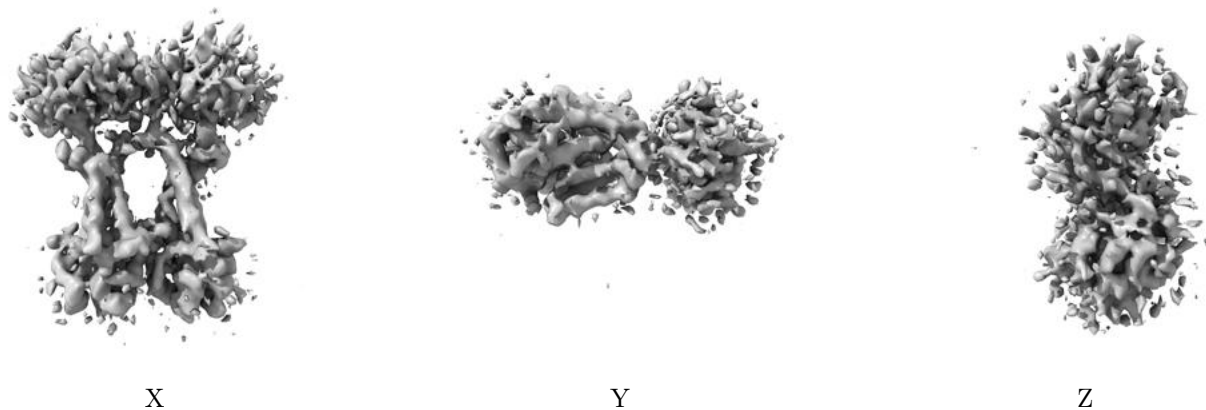


Z

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.

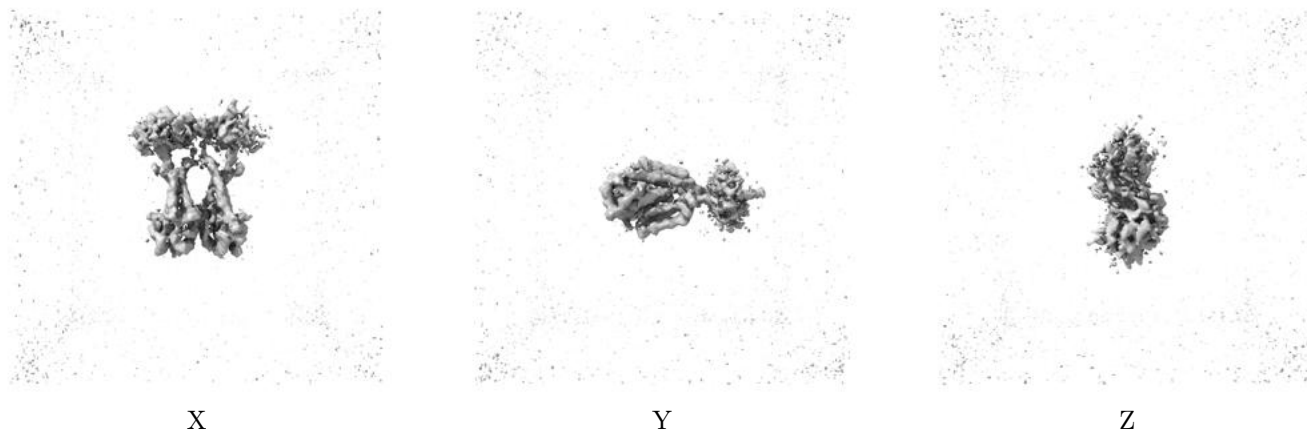
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.17. 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.

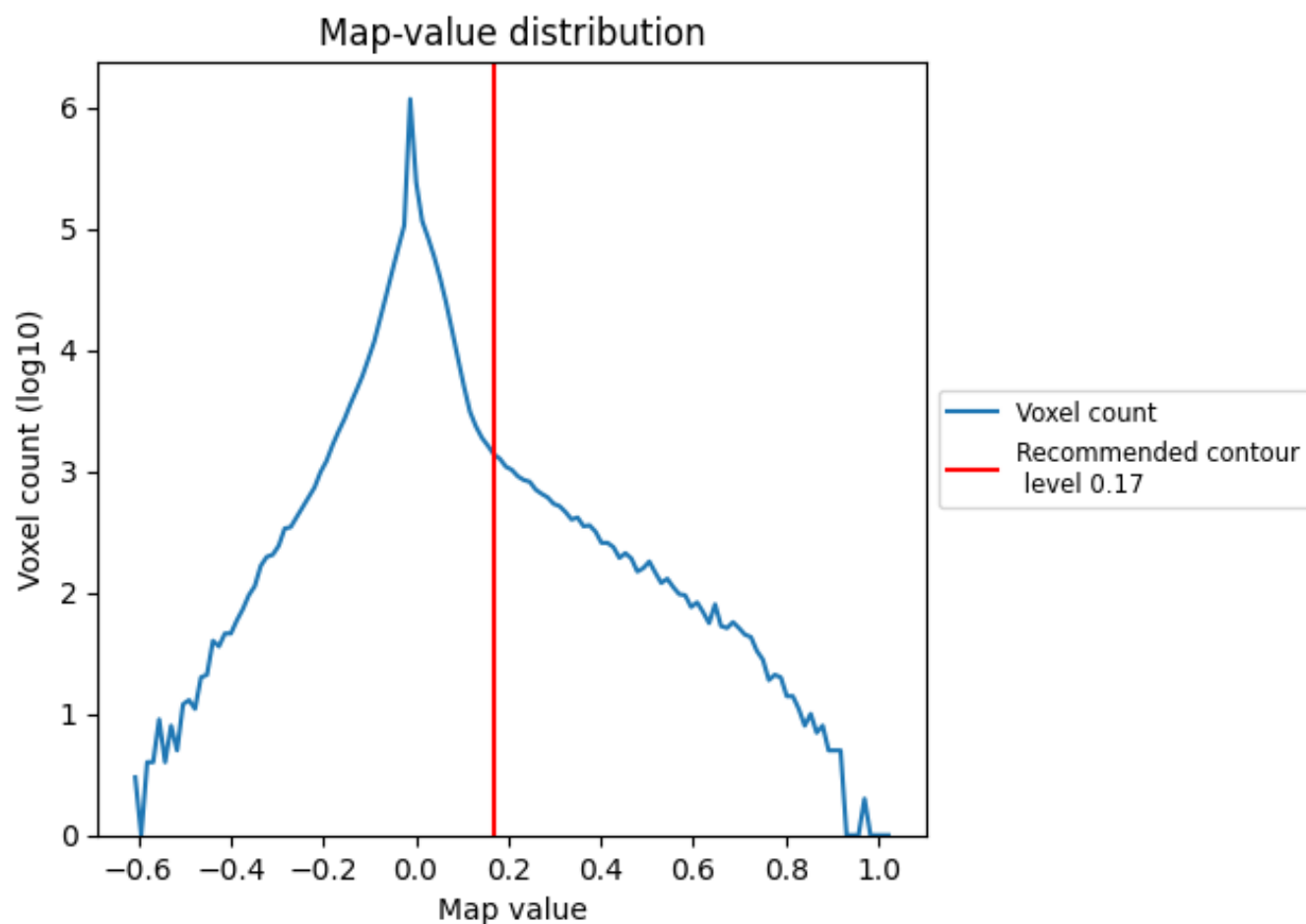
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

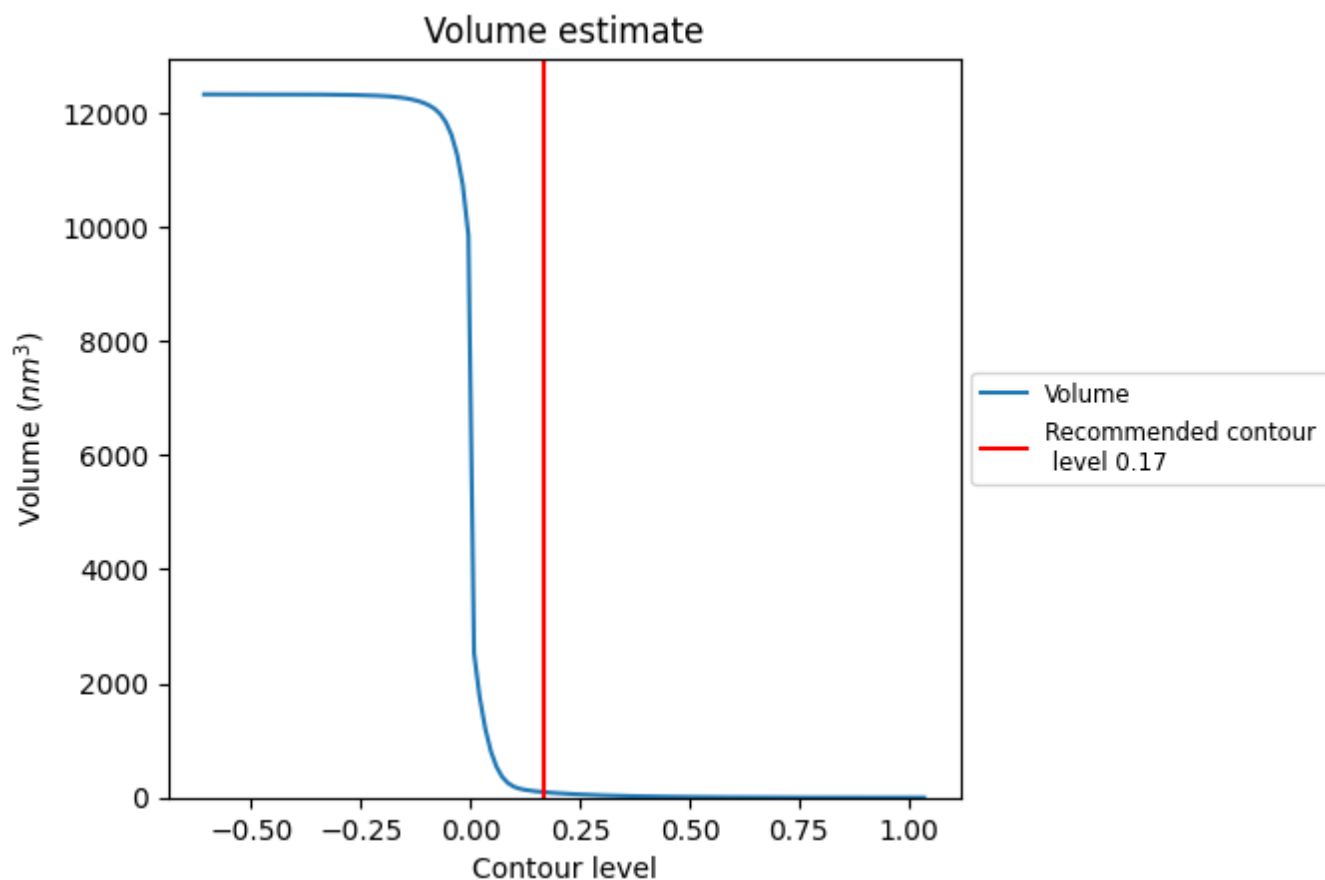
This section contains the results of statistical analysis of the map.

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.

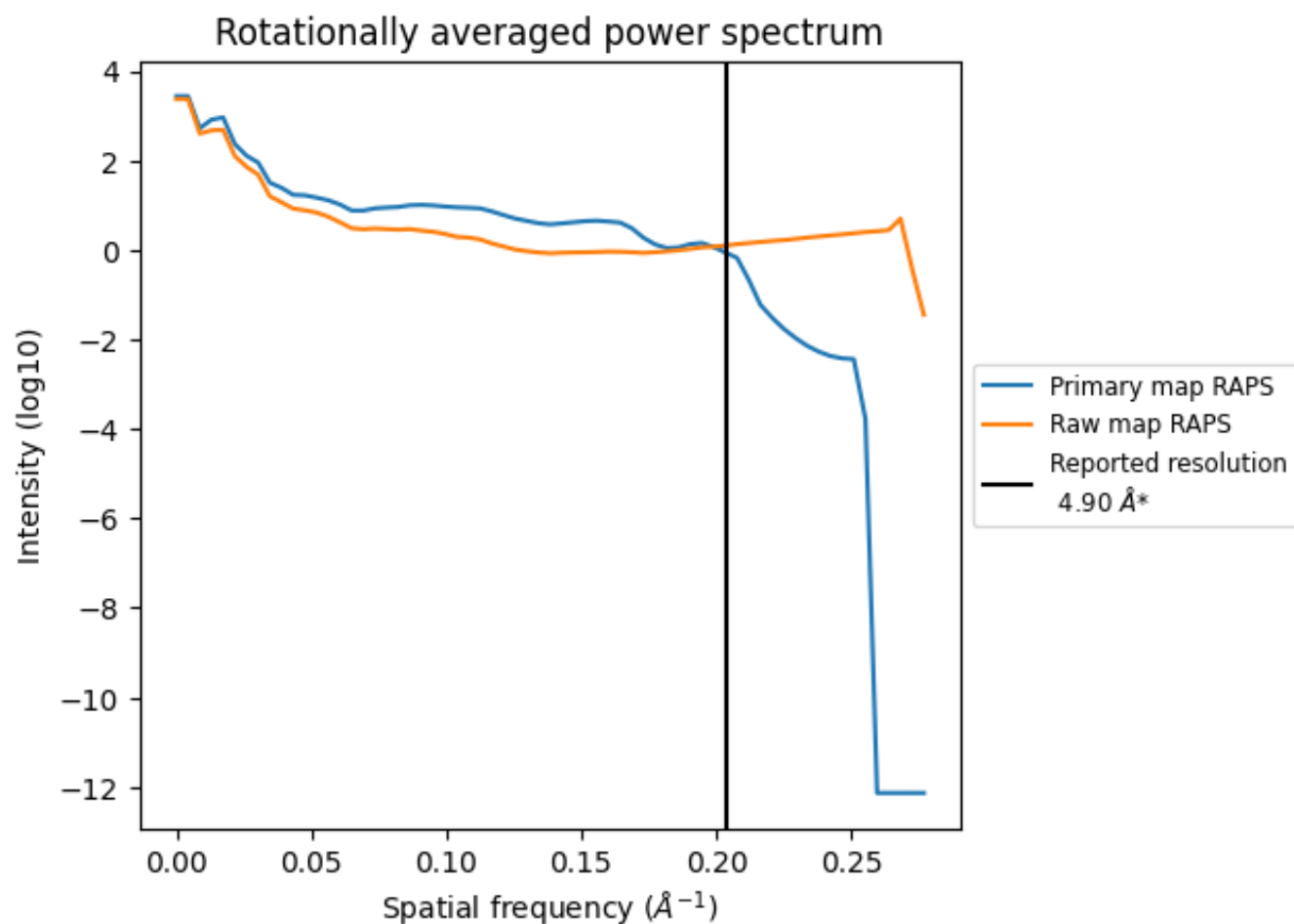
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 94 nm^3 ; this corresponds to an approximate mass of 85 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

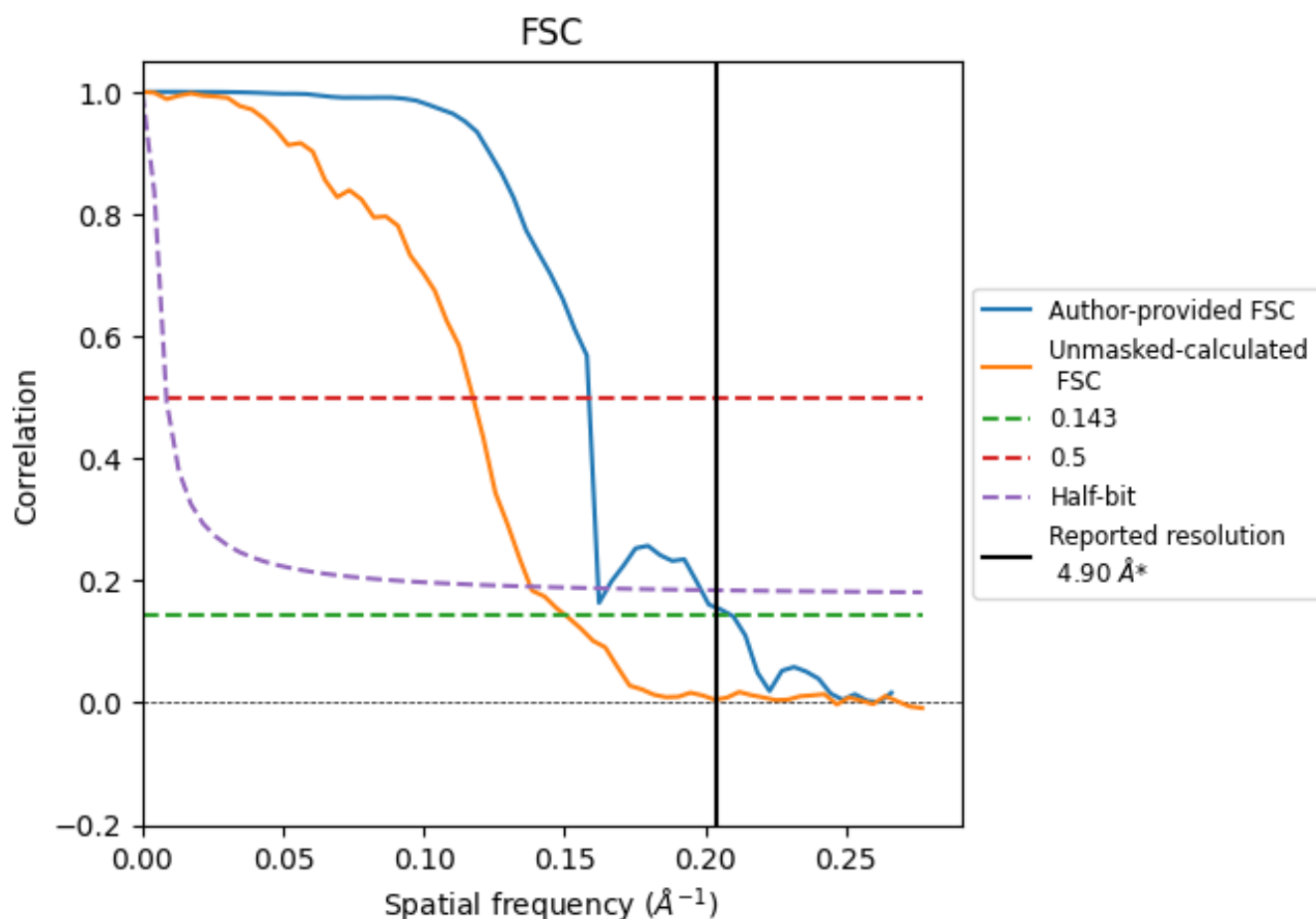


*Reported resolution corresponds to spatial frequency of 0.204 Å⁻¹

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 0.204 Å⁻¹

8.2 Resolution estimates [i](#)

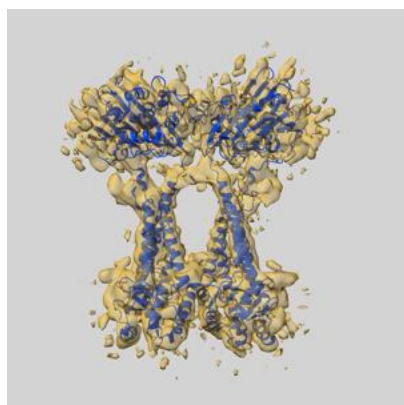
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.90	-	-
Author-provided FSC curve	4.78	6.30	6.17
Unmasked-calculated*	6.65	8.51	7.25

*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 6.65 differs from the reported value 4.9 by more than 10 %

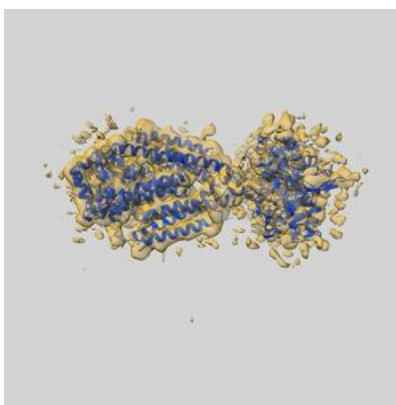
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-28659 and PDB model 8EXQ. Per-residue inclusion information can be found in section [3](#) on page [5](#).

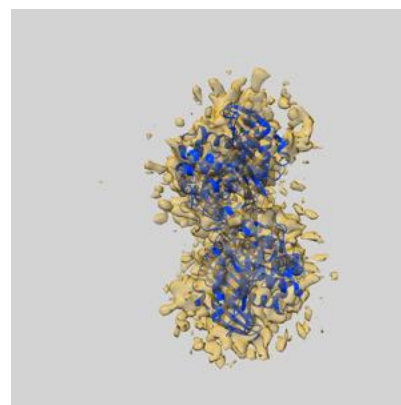
9.1 Map-model overlay [i](#)



X



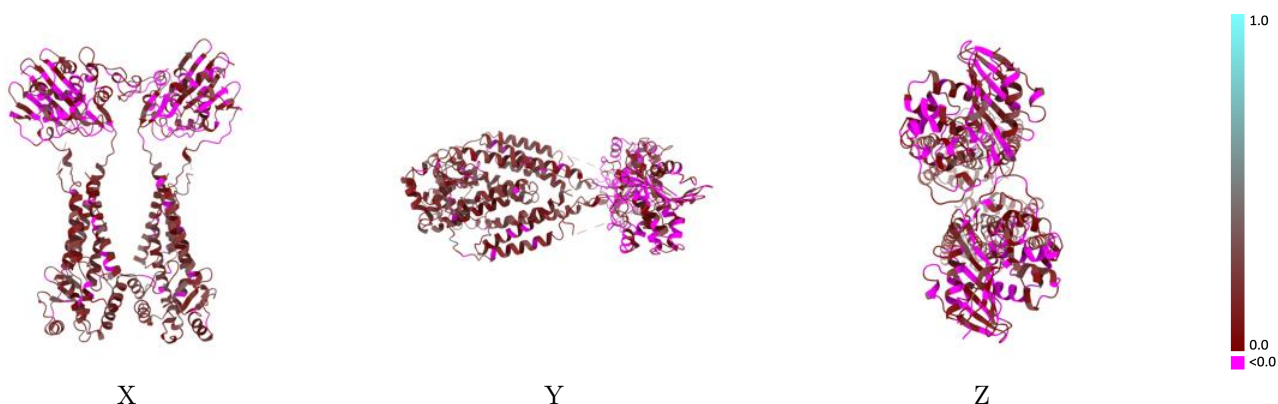
Y



Z

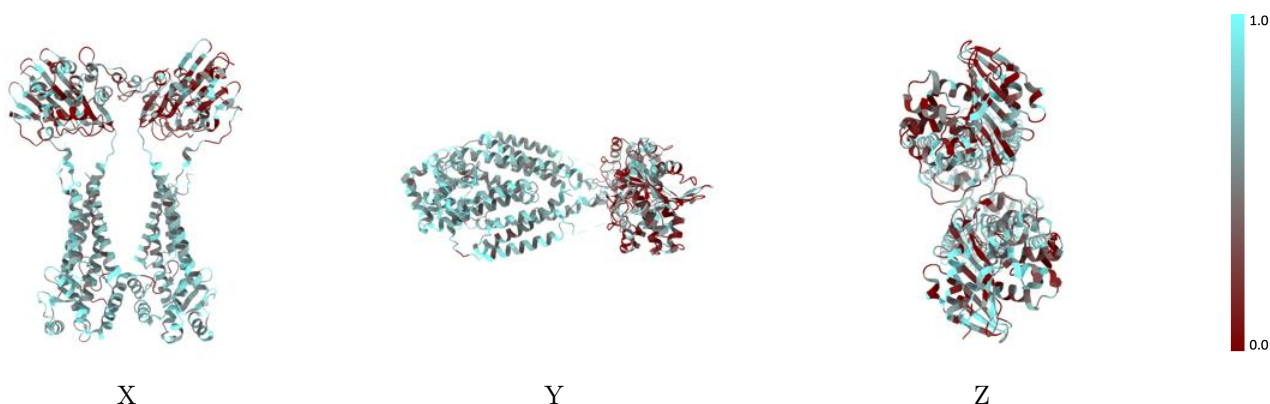
The images above show the 3D surface view of the map at the recommended contour level 0.17 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



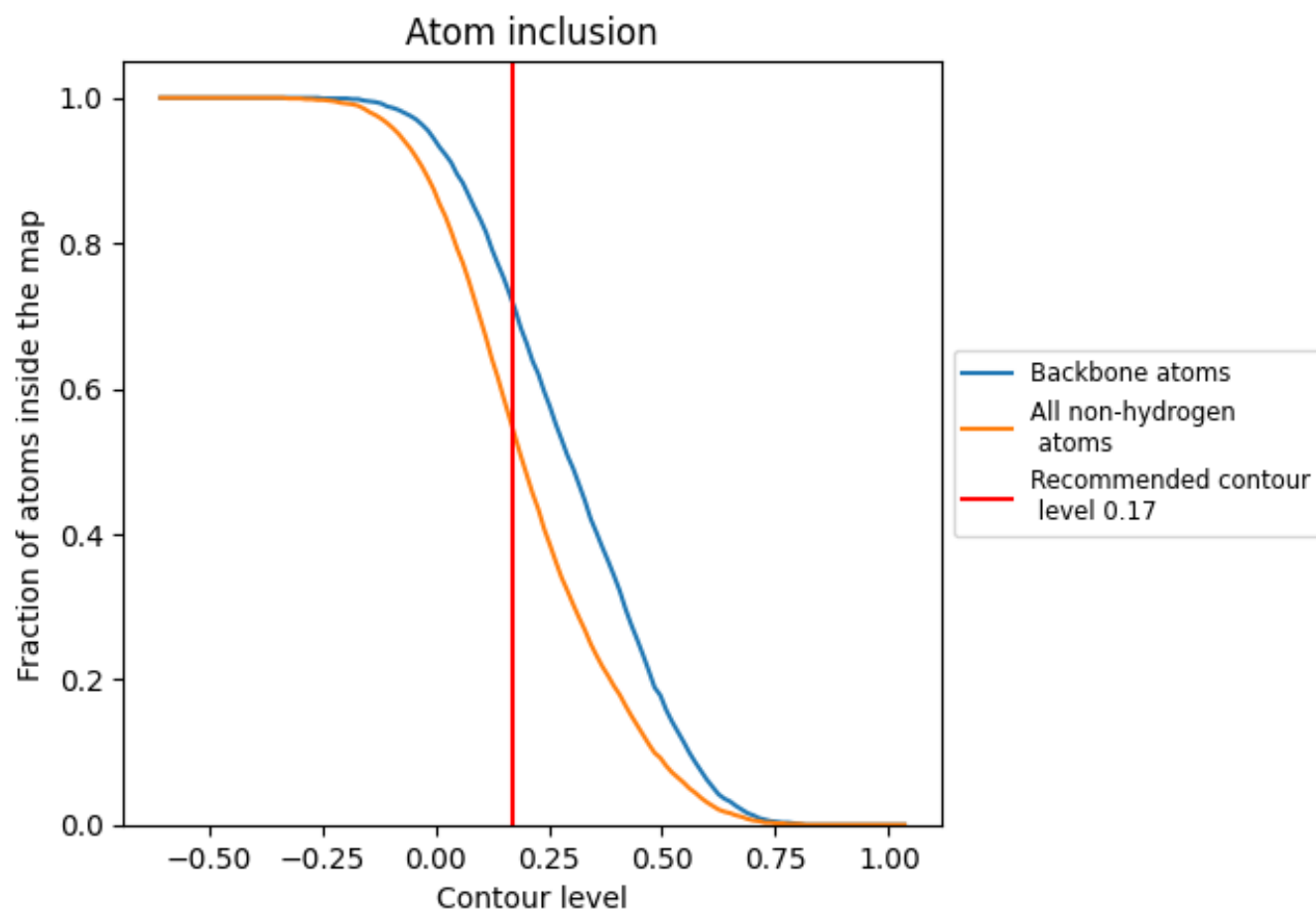
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.17).

9.4 Atom inclusion [i](#)



At the recommended contour level, 72% of all backbone atoms, 55% 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 (0.17) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5460	0.1220
A	0.5230	0.1230
B	0.5690	0.1210

