



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 06:48 AM UTC

PDB ID : 8PHR / pdb_00008phr
EMDB ID : EMD-17672
Title : Middle part of the Borrelia bacteriophage BB1 procapsid, tenfold-symmetrized outer shell
Authors : Rumnieks, J.; Fuzik, T.; Tars, K.
Deposited on : 2023-06-20
Resolution : 2.65 Å(reported)
Based on initial model : .

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:

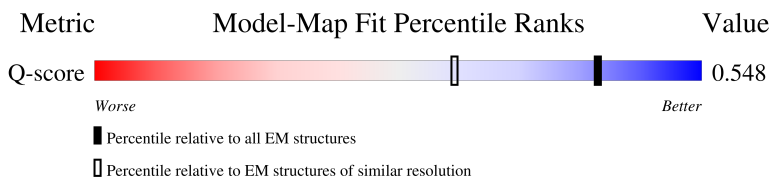
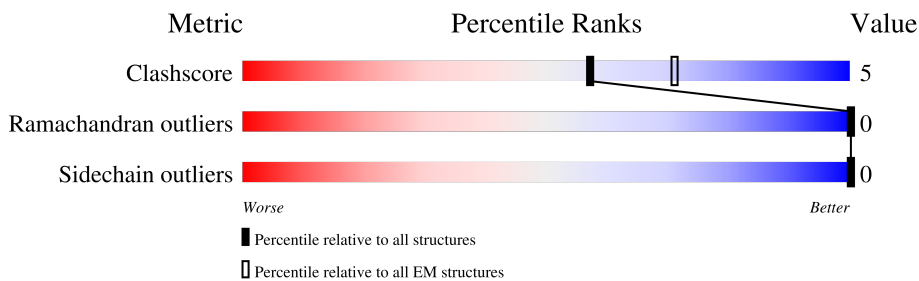
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.65 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.





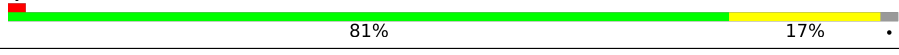




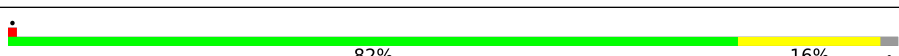
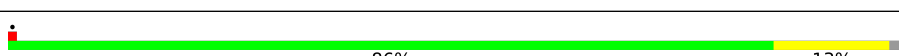

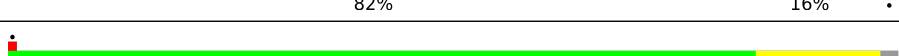
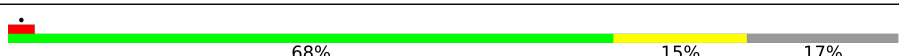
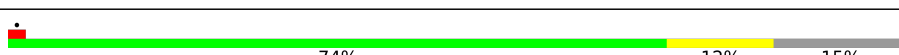

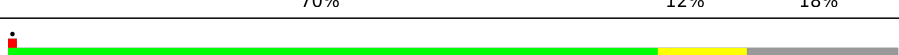
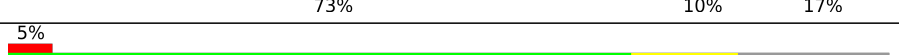
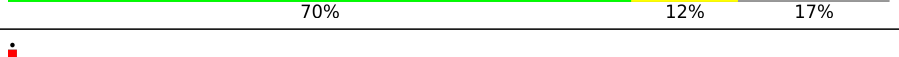




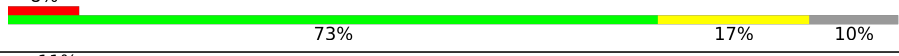



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	9050 (2.15 - 3.15)

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	319	 85% 13%
1	B	319	 85% 13%
1	C	319	 84% 14%
1	J	319	 83% 15%


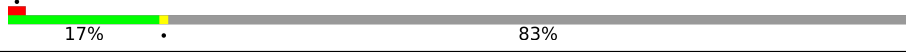




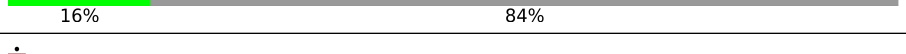
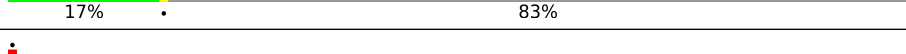
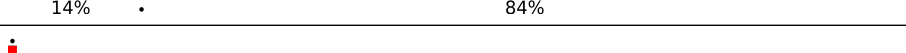

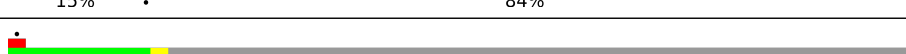


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Mol	Chain	Length	Quality of chain
1	K	319	 83% 15%
1	L	319	 81% 18%
1	S	319	 81% 17%
1	T	319	 86% 13%
1	U	319	 88% 10%
1	b	319	 83% 15%
1	c	319	 83% 15%
1	d	319	 82% 16%
1	k	319	 86% 13%
1	m	319	 82% 16%
1	o	319	 84% 14%
2	D	185	 68% 15% 17%
2	E	185	 74% 12% 15%
2	N	185	 70% 12% 18%
2	O	185	 73% 10% 17%
2	V	185	 5% 70% 12% 17%
2	W	185	 75% 11% 15%
2	f	185	 71% 11% 18%
2	g	185	 74% 10% 17%
3	F	190	 8% 78% 12% 10%
3	M	190	 8% 73% 17% 10%
3	X	190	 11% 78% 12% 10%
3	e	190	 10% 77% 13% 10%
4	G	230	 15% 83%
4	H	230	 15% 84%

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Mol	Chain	Length	Quality of chain
4	I	230	 15% 84%
4	P	230	 17% 83%
4	Q	230	 14% 84%
4	R	230	 16% 84%
4	Y	230	 15% 83%
4	Z	230	 15% 84%
4	a	230	 16% 84%
4	h	230	 17% 83%
4	i	230	 14% 84%
4	j	230	 16% 84%
4	l	230	 15% 84%
4	n	230	 16% 83%
4	p	230	 14% 84%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 57317 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 Major capsid protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	313	2510	1616	404	481	9	0	0
1	B	313	2510	1616	404	481	9	0	0
1	C	313	2510	1616	404	481	9	0	0
1	J	313	2510	1616	404	481	9	0	0
1	K	313	2510	1616	404	481	9	0	0
1	L	313	2510	1616	404	481	9	0	0
1	S	313	2510	1616	404	481	9	0	0
1	T	313	2510	1616	404	481	9	0	0
1	U	313	2510	1616	404	481	9	0	0
1	b	313	2510	1616	404	481	9	0	0
1	c	313	2510	1616	404	481	9	0	0
1	d	313	2510	1616	404	481	9	0	0
1	k	313	2510	1616	404	481	9	0	0
1	m	313	2510	1616	404	481	9	0	0
1	o	313	2510	1616	404	481	9	0	0

- Molecule 2 is a protein called Decorator protein P03.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	153	Total	C	N	O	S	0	0
			1161	730	194	233	4		
2	E	158	Total	C	N	O	S	0	0
			1198	750	202	242	4		
2	N	151	Total	C	N	O	S	0	0
			1148	721	192	231	4		
2	O	154	Total	C	N	O	S	0	0
			1173	739	196	234	4		
2	V	153	Total	C	N	O	S	0	0
			1161	730	194	233	4		
2	W	158	Total	C	N	O	S	0	0
			1198	750	202	242	4		
2	f	151	Total	C	N	O	S	0	0
			1148	721	192	231	4		
2	g	154	Total	C	N	O	S	0	0
			1173	739	196	234	4		

- Molecule 3 is a protein called Decorator protein P05.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	F	171	Total	C	N	O	S	0	0
			1358	865	231	260	2		
3	M	171	Total	C	N	O	S	0	0
			1358	865	231	260	2		
3	X	171	Total	C	N	O	S	0	0
			1358	865	231	260	2		
3	e	171	Total	C	N	O	S	0	0
			1358	865	231	260	2		

- Molecule 4 is a protein called Scaffold protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	G	40	Total	C	N	O	0	0
			345	221	64	60		
4	H	37	Total	C	N	O	0	0
			315	203	55	57		
4	I	37	Total	C	N	O	0	0
			315	203	55	57		
4	P	40	Total	C	N	O	0	0
			345	221	64	60		
4	Q	37	Total	C	N	O	0	0
			315	203	55	57		
4	R	37	Total	C	N	O	0	0
			315	203	55	57		

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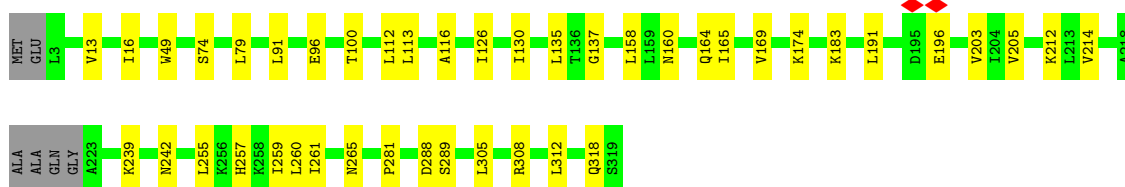
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
4	Y	40	Total 345	C 221	N 64	O 60	0	0
4	Z	37	Total 315	C 203	N 55	O 57	0	0
4	a	37	Total 315	C 203	N 55	O 57	0	0
4	h	40	Total 345	C 221	N 64	O 60	0	0
4	i	37	Total 315	C 203	N 55	O 57	0	0
4	j	37	Total 315	C 203	N 55	O 57	0	0
4	l	37	Total 315	C 203	N 55	O 57	0	0
4	n	40	Total 345	C 221	N 64	O 60	0	0
4	p	37	Total 315	C 203	N 55	O 57	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.

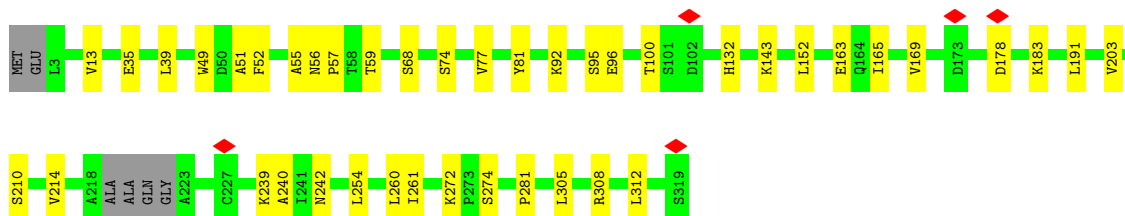
- Molecule 1: Major capsid protein

Chain A: 




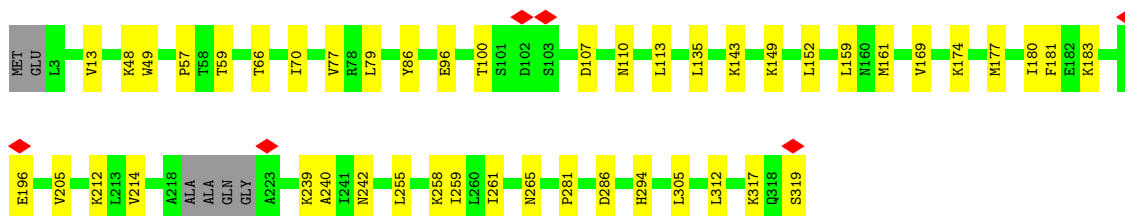
- Molecule 1: Major capsid protein

Chain B: 




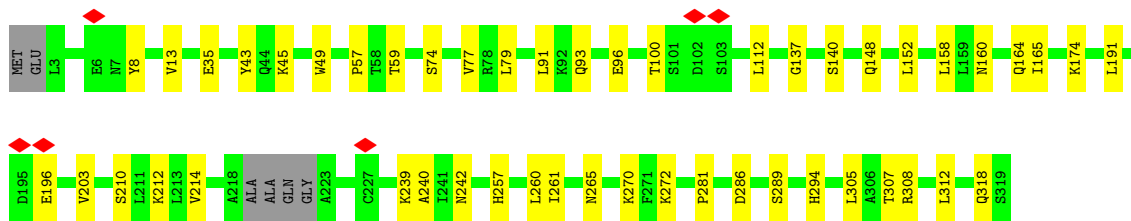
- Molecule 1: Major capsid protein

Chain C: 

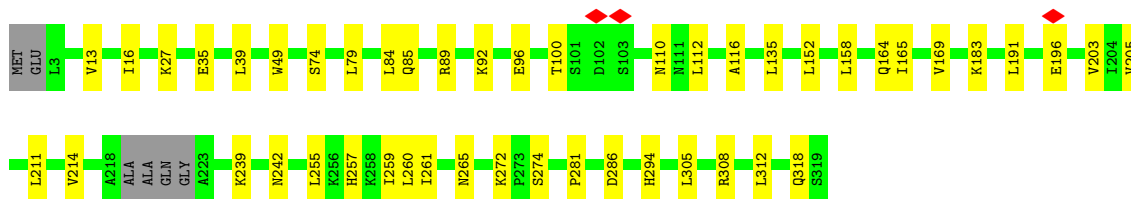
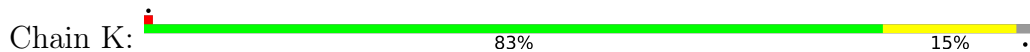


- Molecule 1: Major capsid protein

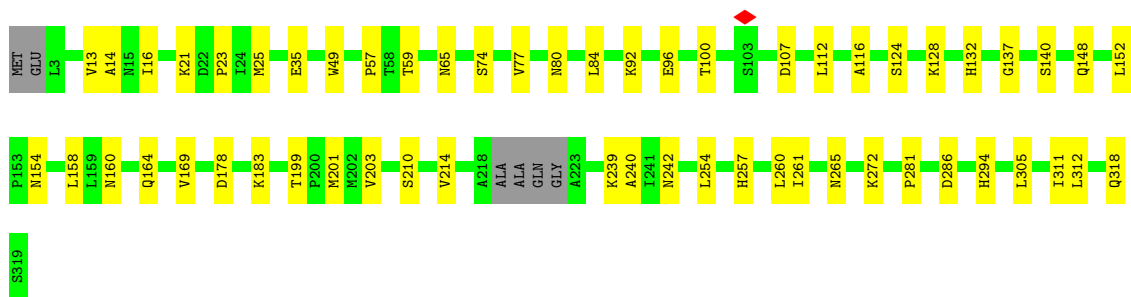
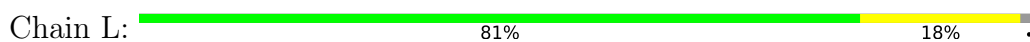
Chain J: 



• Molecule 1: Major capsid protein

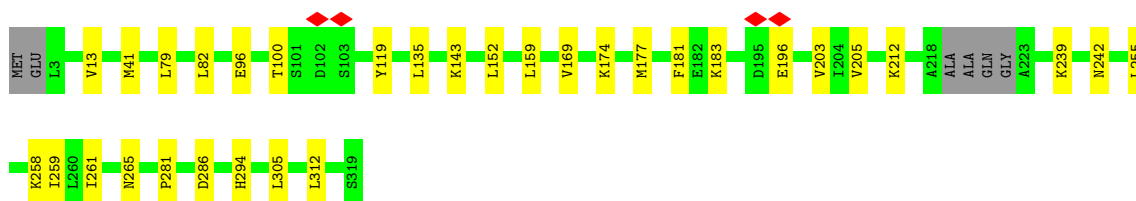
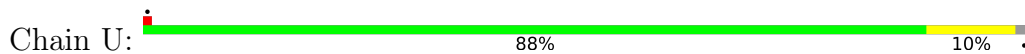


• Molecule 1: Major capsid protein

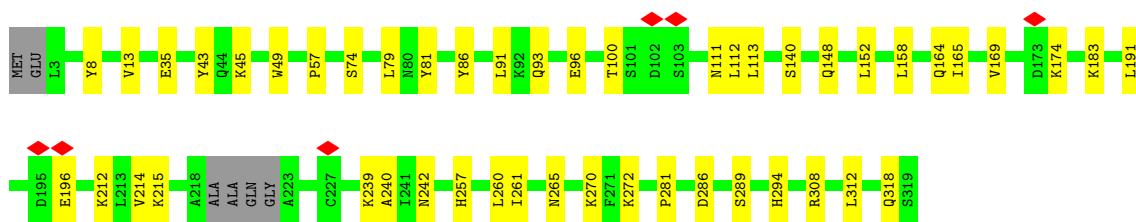
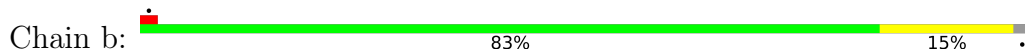




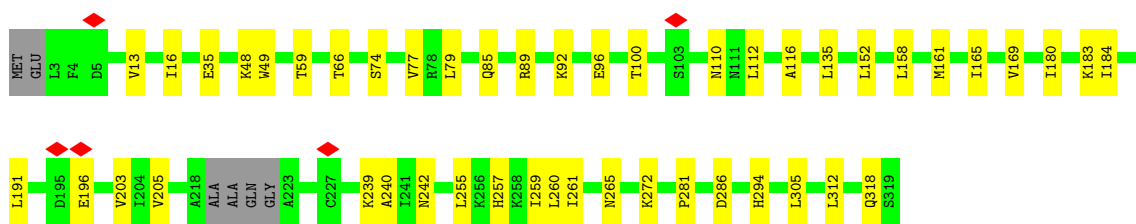
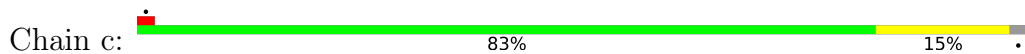
• Molecule 1: Major capsid protein



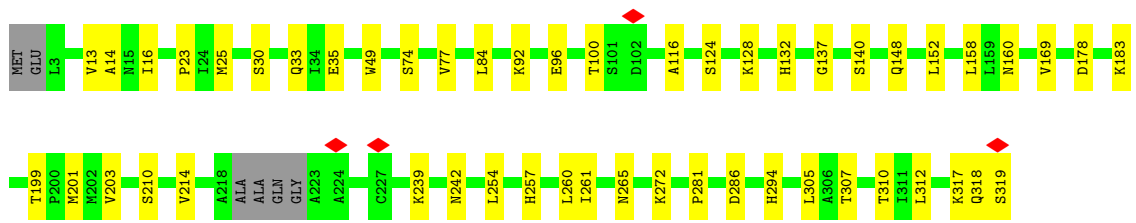
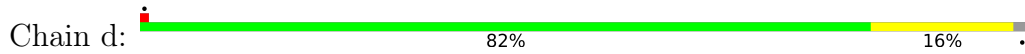
• Molecule 1: Major capsid protein



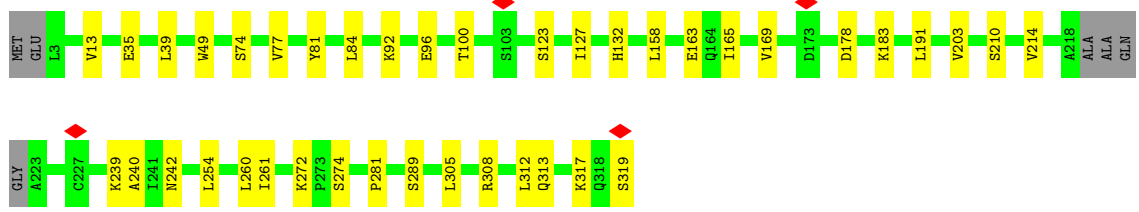
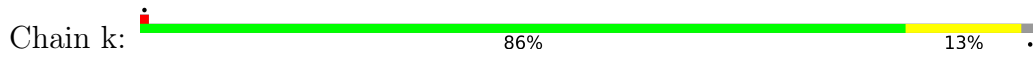
• Molecule 1: Major capsid protein



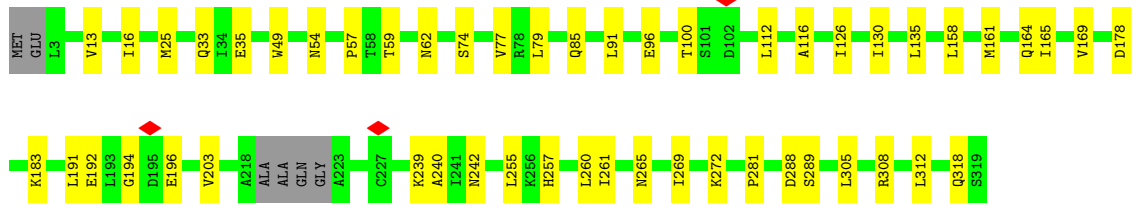
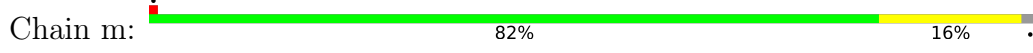
• Molecule 1: Major capsid protein



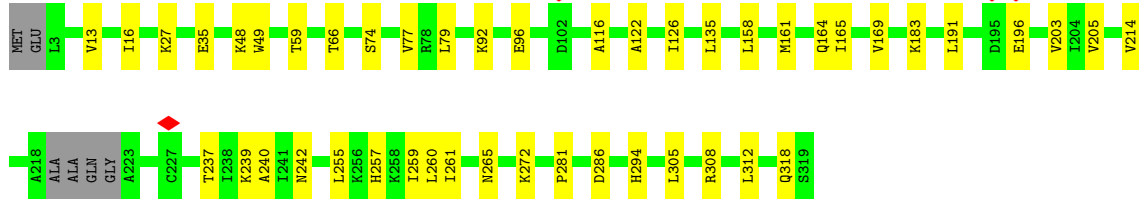
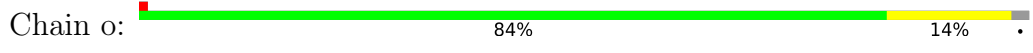
• Molecule 1: Major capsid protein



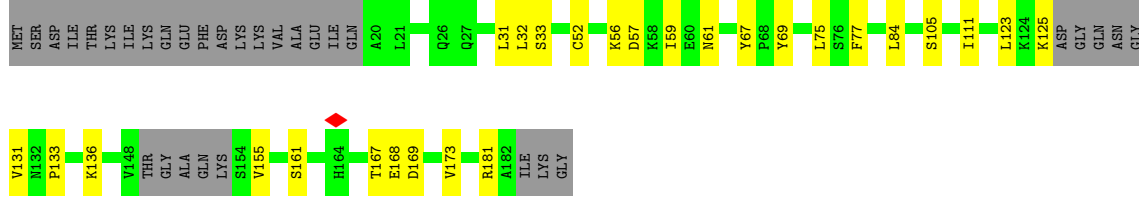
• Molecule 1: Major capsid protein



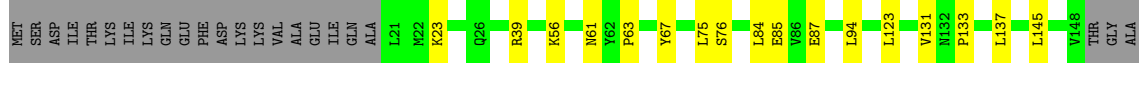
• Molecule 1: Major capsid protein



• Molecule 2: Decorator protein P03



• Molecule 2: Decorator protein P03





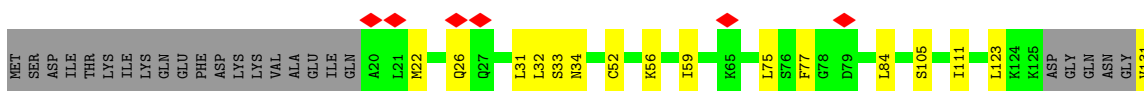
• Molecule 2: Decorator protein P03



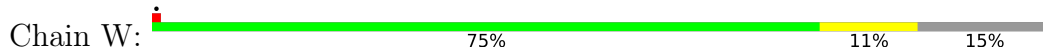
• Molecule 2: Decorator protein P03



• Molecule 2: Decorator protein P03

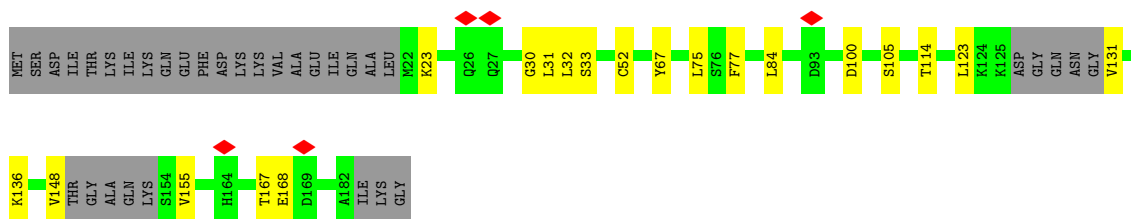


• Molecule 2: Decorator protein P03

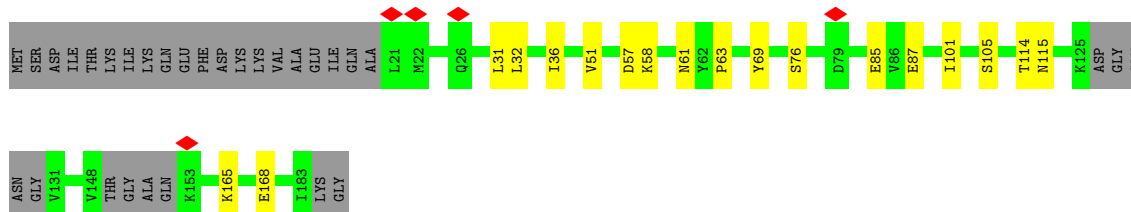


• Molecule 2: Decorator protein P03

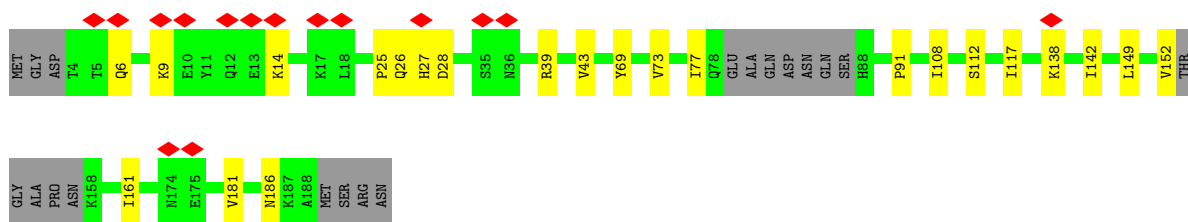
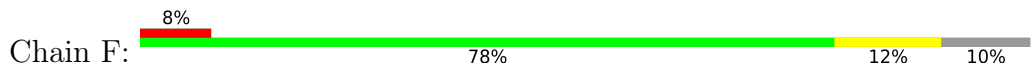




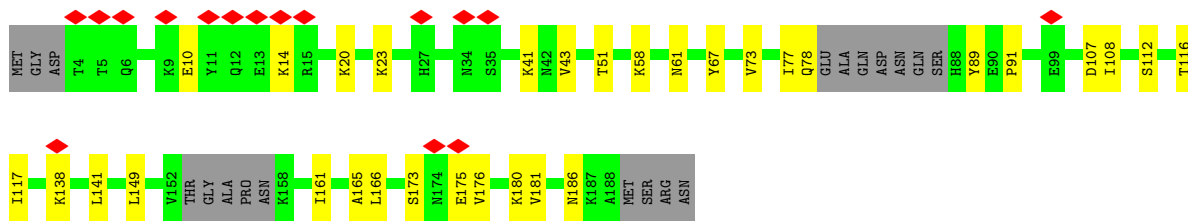
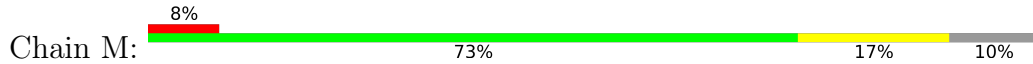
• Molecule 2: Decorator protein P03



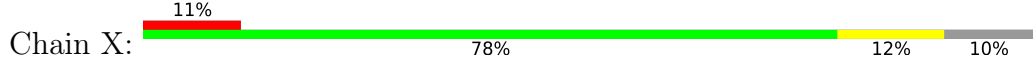
• Molecule 3: Decorator protein P05



• Molecule 3: Decorator protein P05



• Molecule 3: Decorator protein P05



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, D5	Depositor
Number of particles used	29011	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 ($e^-/\text{\AA}^2$)	44	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	165000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.037	Depositor
Minimum map value	-0.019	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.005	Depositor
Map size (Å)	533.50397, 533.50397, 533.50397	wwPDB
Map dimensions	640, 640, 640	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8335999, 0.8335999, 0.8335999	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	0.09	0/2558	0.23	0/3458
1	B	0.08	0/2558	0.23	0/3458
1	C	0.08	0/2558	0.23	0/3458
1	J	0.09	0/2558	0.24	0/3458
1	K	0.09	0/2558	0.23	0/3458
1	L	0.08	0/2558	0.23	0/3458
1	S	0.08	0/2558	0.22	0/3458
1	T	0.09	0/2558	0.23	0/3458
1	U	0.08	0/2558	0.23	0/3458
1	b	0.08	0/2558	0.24	0/3458
1	c	0.08	0/2558	0.23	0/3458
1	d	0.09	0/2558	0.24	0/3458
1	k	0.08	0/2558	0.23	0/3458
1	m	0.08	0/2558	0.22	0/3458
1	o	0.08	0/2558	0.23	0/3458
2	D	0.08	0/1178	0.23	0/1586
2	E	0.08	0/1216	0.22	0/1637
2	N	0.08	0/1165	0.23	0/1568
2	O	0.07	0/1190	0.21	0/1601
2	V	0.09	0/1178	0.22	0/1586
2	W	0.08	0/1216	0.22	0/1637
2	f	0.08	0/1165	0.22	0/1568
2	g	0.07	0/1190	0.23	0/1601
3	F	0.08	0/1381	0.23	0/1855
3	M	0.09	0/1381	0.24	0/1855
3	X	0.08	0/1381	0.23	0/1855
3	e	0.08	0/1381	0.24	0/1855
4	G	0.08	0/352	0.19	0/472
4	H	0.08	0/321	0.19	0/432
4	I	0.07	0/321	0.15	0/432
4	P	0.07	0/352	0.17	0/472
4	Q	0.08	0/321	0.17	0/432
4	R	0.07	0/321	0.17	0/432
4	Y	0.08	0/352	0.19	0/472

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
4	Z	0.08	0/321	0.18	0/432
4	a	0.07	0/321	0.15	0/432
4	h	0.07	0/352	0.16	0/472
4	i	0.08	0/321	0.17	0/432
4	j	0.07	0/321	0.16	0/432
4	l	0.07	0/321	0.17	0/432
4	n	0.08	0/352	0.19	0/472
4	p	0.07	0/321	0.18	0/432
All	All	0.08	0/58362	0.22	0/78754

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	A	2510	0	2547	24	0
1	B	2510	0	2547	28	0
1	C	2510	0	2547	33	0
1	J	2510	0	2547	31	0
1	K	2510	0	2547	29	0
1	L	2510	0	2547	37	0
1	S	2510	0	2547	38	0
1	T	2510	0	2547	25	0
1	U	2510	0	2547	18	0
1	b	2510	0	2547	32	0
1	c	2510	0	2547	32	0
1	d	2510	0	2547	30	0
1	k	2510	0	2547	25	0
1	m	2510	0	2547	33	0
1	o	2510	0	2547	29	0
2	D	1161	0	1139	20	0
2	E	1198	0	1172	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	N	1148	0	1123	14	0
2	O	1173	0	1158	15	0
2	V	1161	0	1139	17	0
2	W	1198	0	1172	15	0
2	f	1148	0	1123	15	0
2	g	1173	0	1158	14	0
3	F	1358	0	1361	15	0
3	M	1358	0	1361	22	0
3	X	1358	0	1361	16	0
3	e	1358	0	1361	15	0
4	G	345	0	345	3	0
4	H	315	0	312	1	0
4	I	315	0	312	1	0
4	P	345	0	345	1	0
4	Q	315	0	312	3	0
4	R	315	0	312	1	0
4	Y	345	0	345	3	0
4	Z	315	0	312	1	0
4	a	315	0	312	0	0
4	h	345	0	345	1	0
4	i	315	0	312	2	0
4	j	315	0	312	1	0
4	l	315	0	312	1	0
4	n	345	0	345	2	0
4	p	315	0	312	3	0
All	All	57317	0	57678	553	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 5.

The worst 5 of 553 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:A:165:ILE:HD13	1:A:191:LEU:HD23	1.70	0.73
1:c:165:ILE:HD13	1:c:191:LEU:HD23	1.70	0.72
1:S:165:ILE:HD13	1:S:191:LEU:HD23	1.72	0.72
1:o:165:ILE:HD13	1:o:191:LEU:HD23	1.70	0.71
1:K:165:ILE:HD13	1:K:191:LEU:HD23	1.71	0.71

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	309/319 (97%)	305 (99%)	4 (1%)	0	100	100
1	B	309/319 (97%)	306 (99%)	3 (1%)	0	100	100
1	C	309/319 (97%)	306 (99%)	3 (1%)	0	100	100
1	J	309/319 (97%)	302 (98%)	7 (2%)	0	100	100
1	K	309/319 (97%)	307 (99%)	2 (1%)	0	100	100
1	L	309/319 (97%)	305 (99%)	4 (1%)	0	100	100
1	S	309/319 (97%)	305 (99%)	4 (1%)	0	100	100
1	T	309/319 (97%)	307 (99%)	2 (1%)	0	100	100
1	U	309/319 (97%)	305 (99%)	4 (1%)	0	100	100
1	b	309/319 (97%)	301 (97%)	8 (3%)	0	100	100
1	c	309/319 (97%)	305 (99%)	4 (1%)	0	100	100
1	d	309/319 (97%)	305 (99%)	4 (1%)	0	100	100
1	k	309/319 (97%)	307 (99%)	2 (1%)	0	100	100
1	m	309/319 (97%)	305 (99%)	4 (1%)	0	100	100
1	o	309/319 (97%)	305 (99%)	4 (1%)	0	100	100
2	D	147/185 (80%)	145 (99%)	2 (1%)	0	100	100
2	E	154/185 (83%)	153 (99%)	1 (1%)	0	100	100
2	N	145/185 (78%)	144 (99%)	1 (1%)	0	100	100
2	O	148/185 (80%)	146 (99%)	2 (1%)	0	100	100
2	V	147/185 (80%)	146 (99%)	1 (1%)	0	100	100
2	W	154/185 (83%)	152 (99%)	2 (1%)	0	100	100
2	f	145/185 (78%)	145 (100%)	0	0	100	100
2	g	148/185 (80%)	147 (99%)	1 (1%)	0	100	100
3	F	165/190 (87%)	164 (99%)	1 (1%)	0	100	100
3	M	165/190 (87%)	163 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	X	165/190 (87%)	163 (99%)	2 (1%)	0	100	100
3	e	165/190 (87%)	162 (98%)	3 (2%)	0	100	100
4	G	38/230 (16%)	38 (100%)	0	0	100	100
4	H	35/230 (15%)	35 (100%)	0	0	100	100
4	I	35/230 (15%)	35 (100%)	0	0	100	100
4	P	38/230 (16%)	38 (100%)	0	0	100	100
4	Q	35/230 (15%)	35 (100%)	0	0	100	100
4	R	35/230 (15%)	35 (100%)	0	0	100	100
4	Y	38/230 (16%)	38 (100%)	0	0	100	100
4	Z	35/230 (15%)	35 (100%)	0	0	100	100
4	a	35/230 (15%)	35 (100%)	0	0	100	100
4	h	38/230 (16%)	38 (100%)	0	0	100	100
4	i	35/230 (15%)	35 (100%)	0	0	100	100
4	j	35/230 (15%)	35 (100%)	0	0	100	100
4	l	35/230 (15%)	35 (100%)	0	0	100	100
4	n	38/230 (16%)	38 (100%)	0	0	100	100
4	p	35/230 (15%)	35 (100%)	0	0	100	100
All	All	7023/10475 (67%)	6946 (99%)	77 (1%)	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	283/286 (99%)	283 (100%)	0	100	100
1	B	283/286 (99%)	283 (100%)	0	100	100
1	C	283/286 (99%)	283 (100%)	0	100	100
1	J	283/286 (99%)	283 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	K	283/286 (99%)	283 (100%)	0	100	100
1	L	283/286 (99%)	283 (100%)	0	100	100
1	S	283/286 (99%)	283 (100%)	0	100	100
1	T	283/286 (99%)	283 (100%)	0	100	100
1	U	283/286 (99%)	283 (100%)	0	100	100
1	b	283/286 (99%)	283 (100%)	0	100	100
1	c	283/286 (99%)	283 (100%)	0	100	100
1	d	283/286 (99%)	283 (100%)	0	100	100
1	k	283/286 (99%)	283 (100%)	0	100	100
1	m	283/286 (99%)	283 (100%)	0	100	100
1	o	283/286 (99%)	283 (100%)	0	100	100
2	D	128/154 (83%)	128 (100%)	0	100	100
2	E	132/154 (86%)	132 (100%)	0	100	100
2	N	127/154 (82%)	127 (100%)	0	100	100
2	O	130/154 (84%)	130 (100%)	0	100	100
2	V	128/154 (83%)	128 (100%)	0	100	100
2	W	132/154 (86%)	132 (100%)	0	100	100
2	f	127/154 (82%)	127 (100%)	0	100	100
2	g	130/154 (84%)	130 (100%)	0	100	100
3	F	148/163 (91%)	148 (100%)	0	100	100
3	M	148/163 (91%)	148 (100%)	0	100	100
3	X	148/163 (91%)	148 (100%)	0	100	100
3	e	148/163 (91%)	148 (100%)	0	100	100
4	G	38/210 (18%)	38 (100%)	0	100	100
4	H	35/210 (17%)	35 (100%)	0	100	100
4	I	35/210 (17%)	35 (100%)	0	100	100
4	P	38/210 (18%)	38 (100%)	0	100	100
4	Q	35/210 (17%)	35 (100%)	0	100	100
4	R	35/210 (17%)	35 (100%)	0	100	100
4	Y	38/210 (18%)	38 (100%)	0	100	100
4	Z	35/210 (17%)	35 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	a	35/210 (17%)	35 (100%)	0	100	100
4	h	38/210 (18%)	38 (100%)	0	100	100
4	i	35/210 (17%)	35 (100%)	0	100	100
4	j	35/210 (17%)	35 (100%)	0	100	100
4	l	35/210 (17%)	35 (100%)	0	100	100
4	n	38/210 (18%)	38 (100%)	0	100	100
4	p	35/210 (17%)	35 (100%)	0	100	100
All	All	6411/9324 (69%)	6411 (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. 5 of 65 such sidechains are listed below:

Mol	Chain	Res	Type
1	m	117	GLN
1	m	236	GLN
1	L	315	ASN
1	L	111	ASN
1	m	315	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](#)

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.

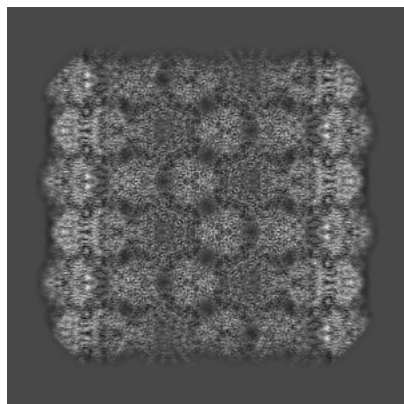
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-17672. These allow visual inspection of the internal detail of the map and identification of artifacts.

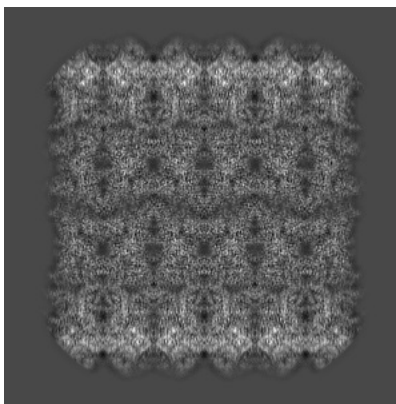
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

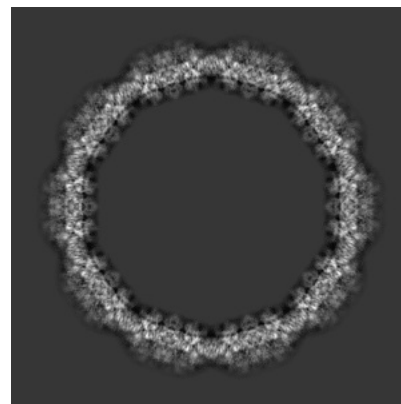
6.1.1 Primary map



X

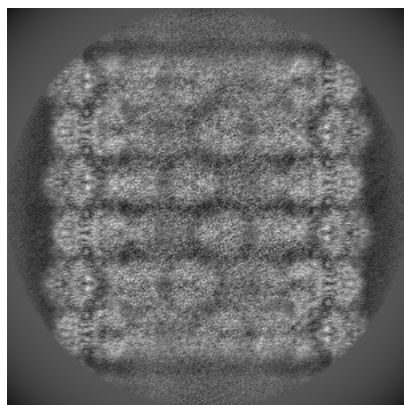


Y

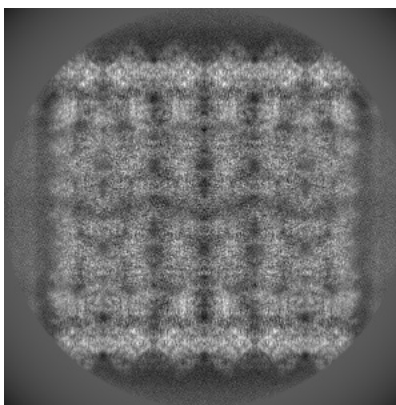


Z

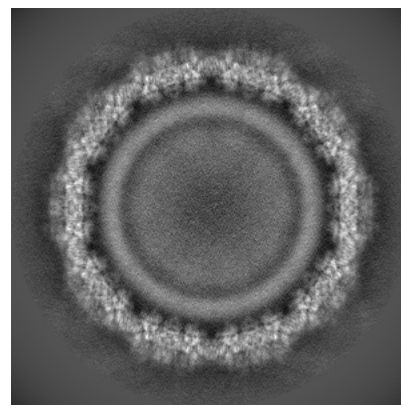
6.1.2 Raw map



X



Y

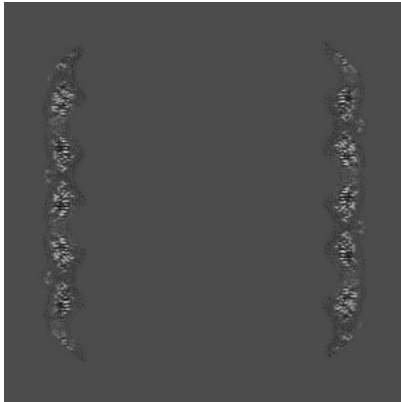


Z

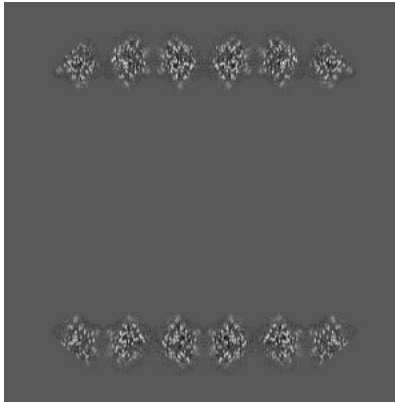
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

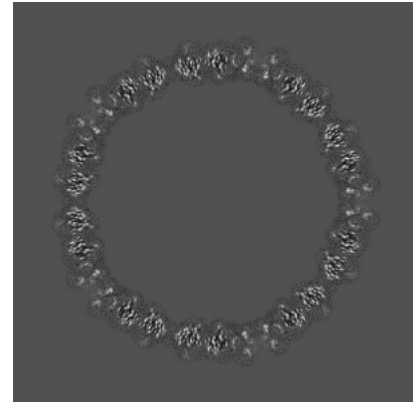
6.2.1 Primary map



X Index: 320

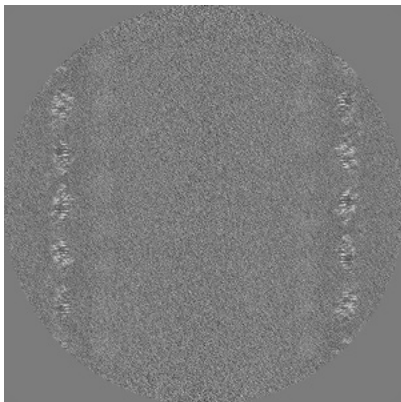


Y Index: 320

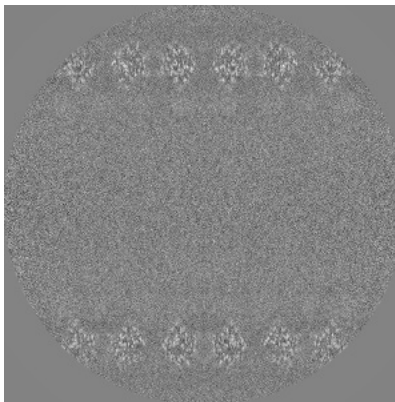


Z Index: 320

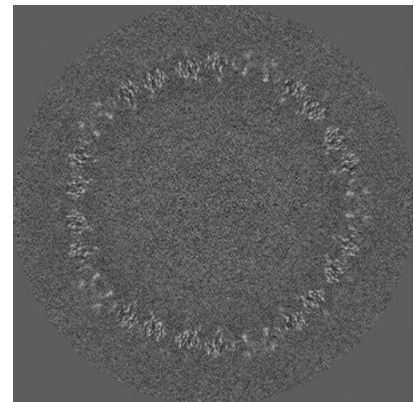
6.2.2 Raw map



X Index: 320



Y Index: 320

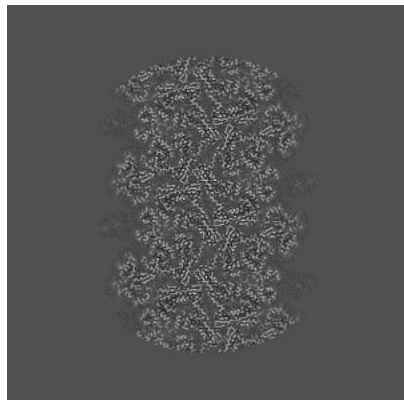


Z Index: 320

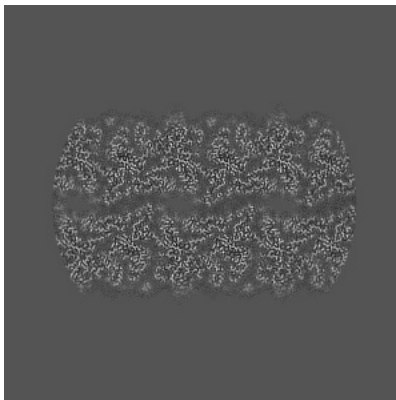
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

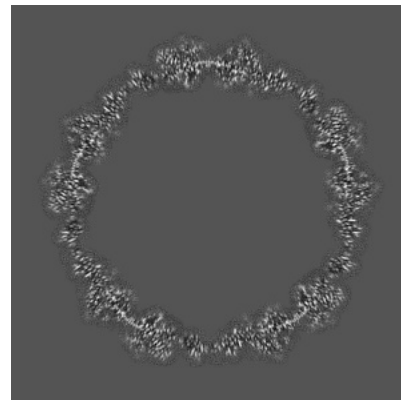
6.3.1 Primary map



X Index: 536

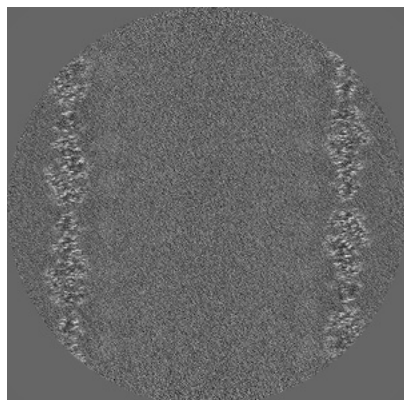


Y Index: 110

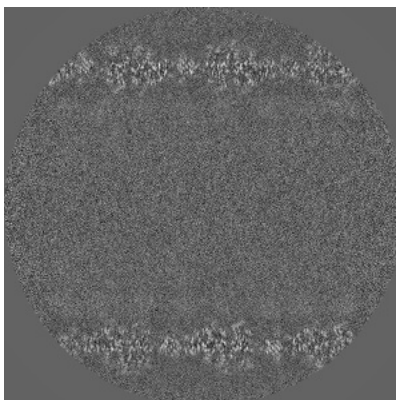


Z Index: 295

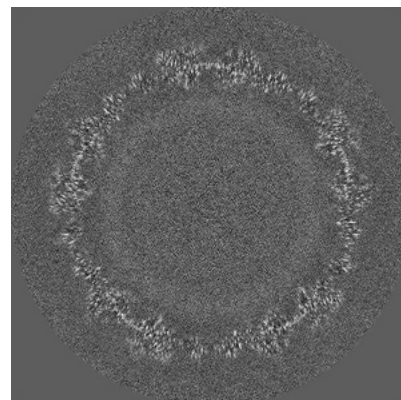
6.3.2 Raw map



X Index: 350



Y Index: 286

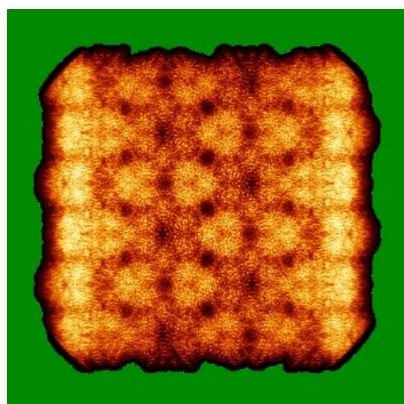


Z Index: 295

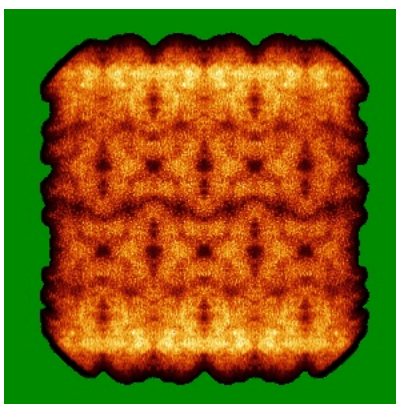
The images above show the largest variance slices of the map in three orthogonal directions.

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

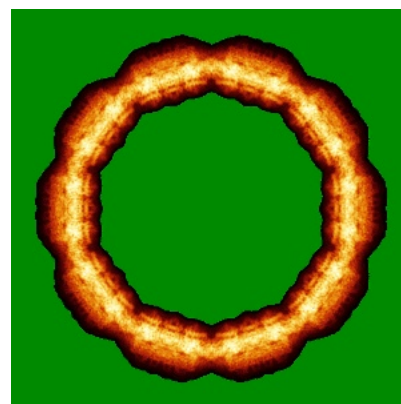
6.4.1 Primary map



X

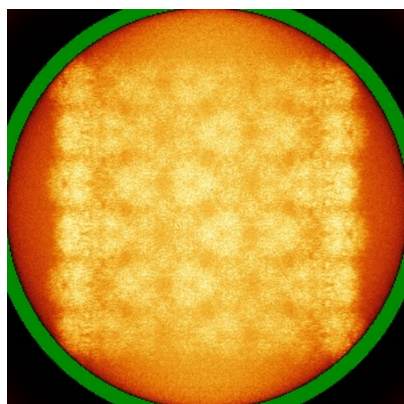


Y

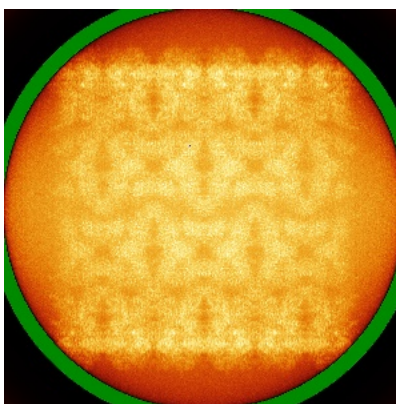


Z

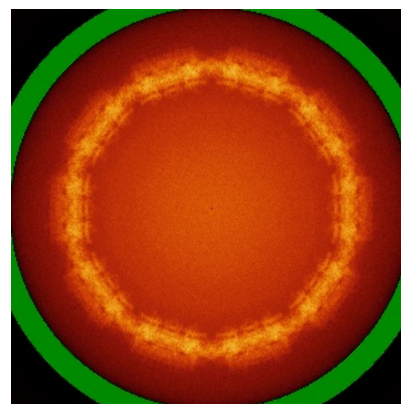
6.4.2 Raw map



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](#)

This section was not generated.

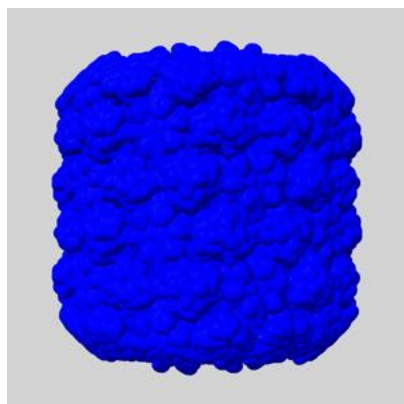
6.6 Mask visualisation [i](#)

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

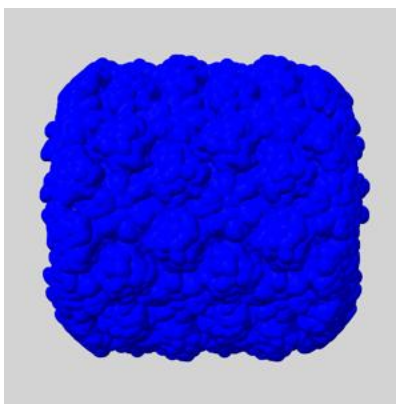
A mask typically either:

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

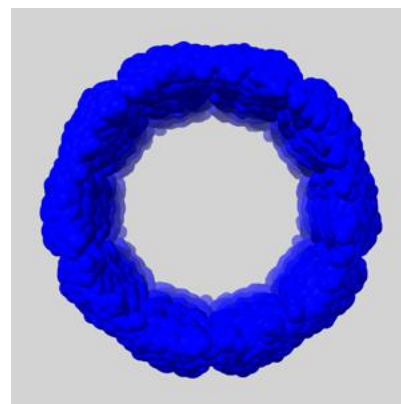
6.6.1 emd_17672_msk_1.map [i](#)



X



Y

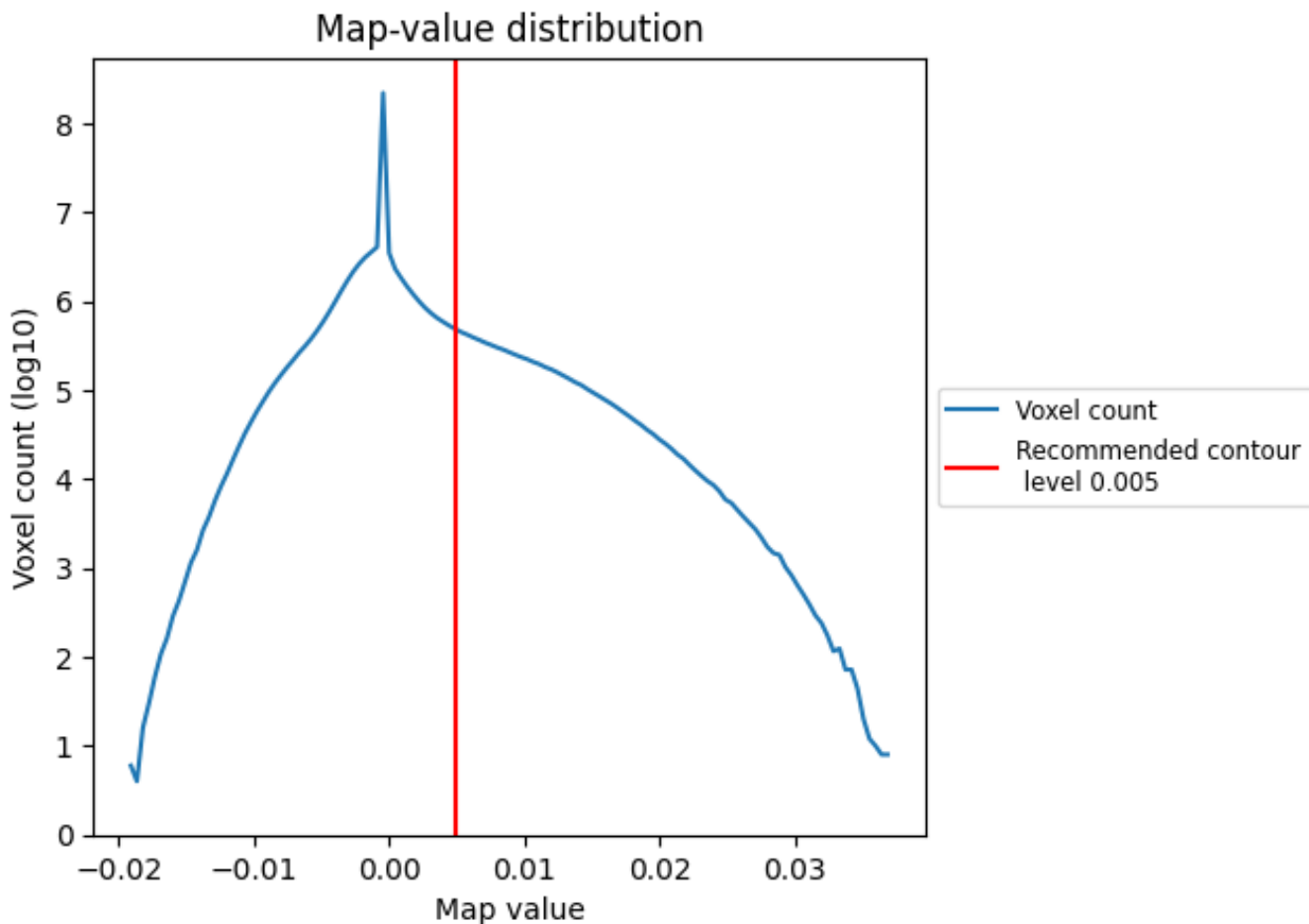


Z

7 Map analysis [i](#)

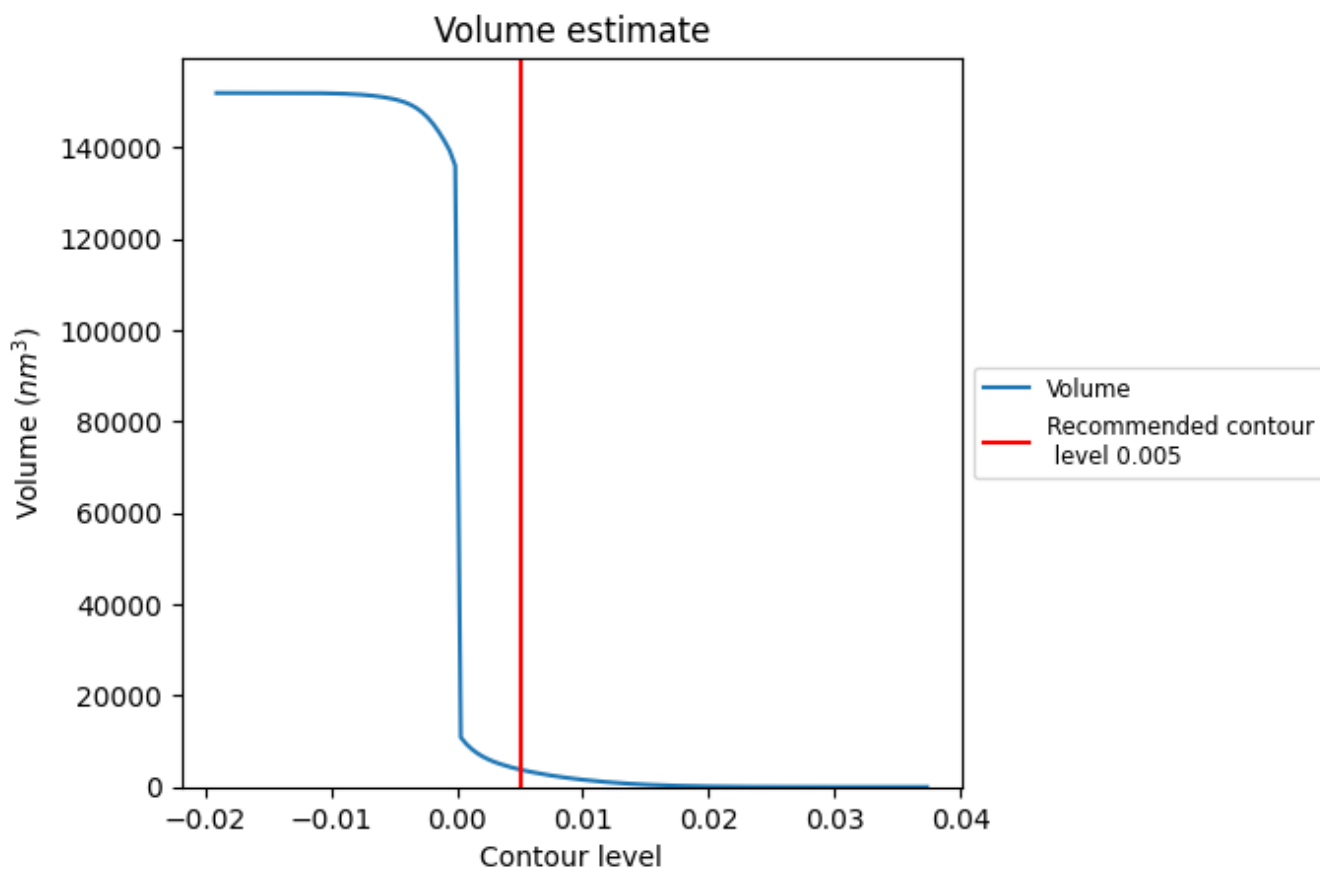
This section contains the results of statistical analysis of the map.

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.

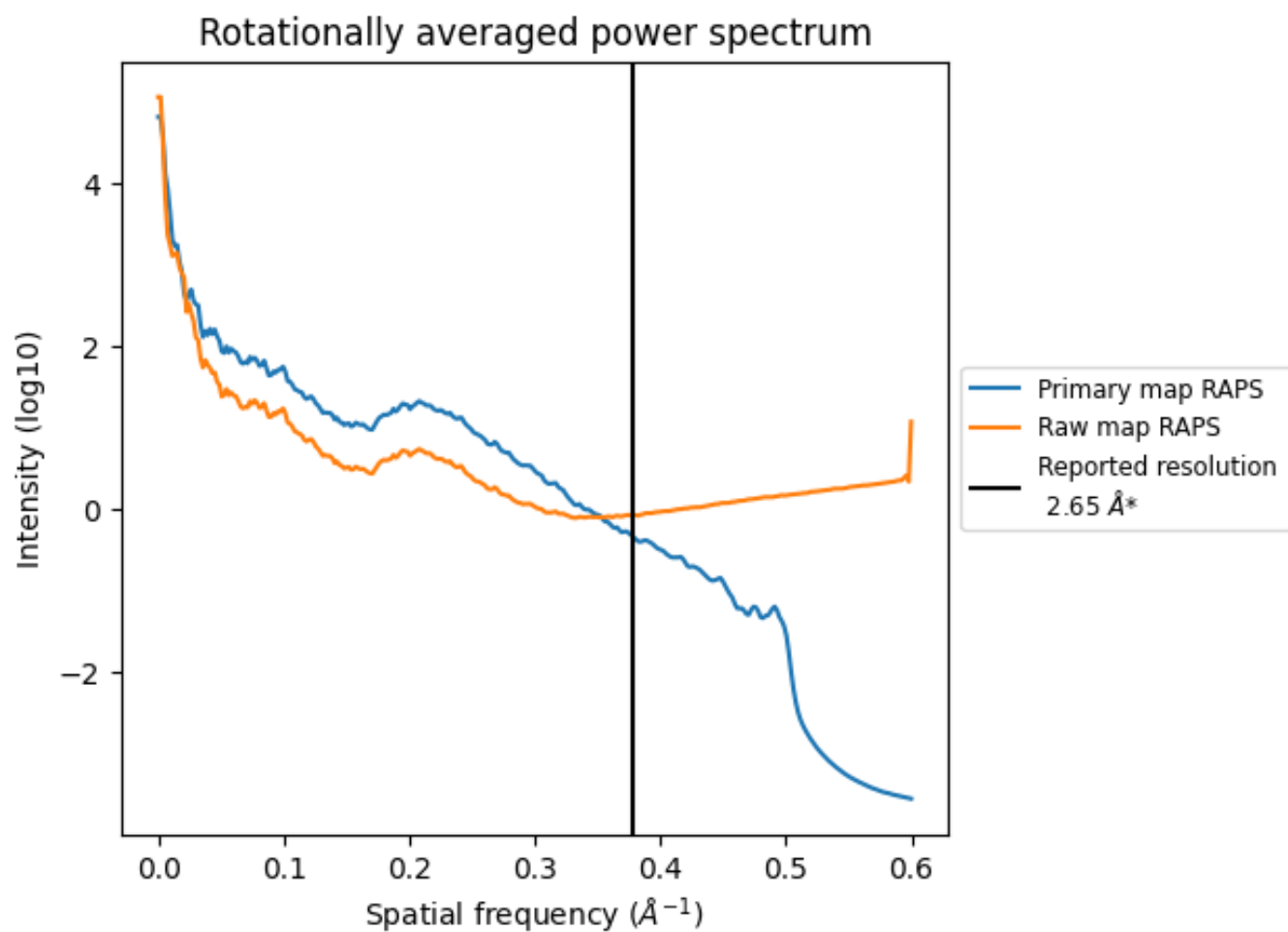
7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 3826 nm^3 ; this corresponds to an approximate mass of 3456 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

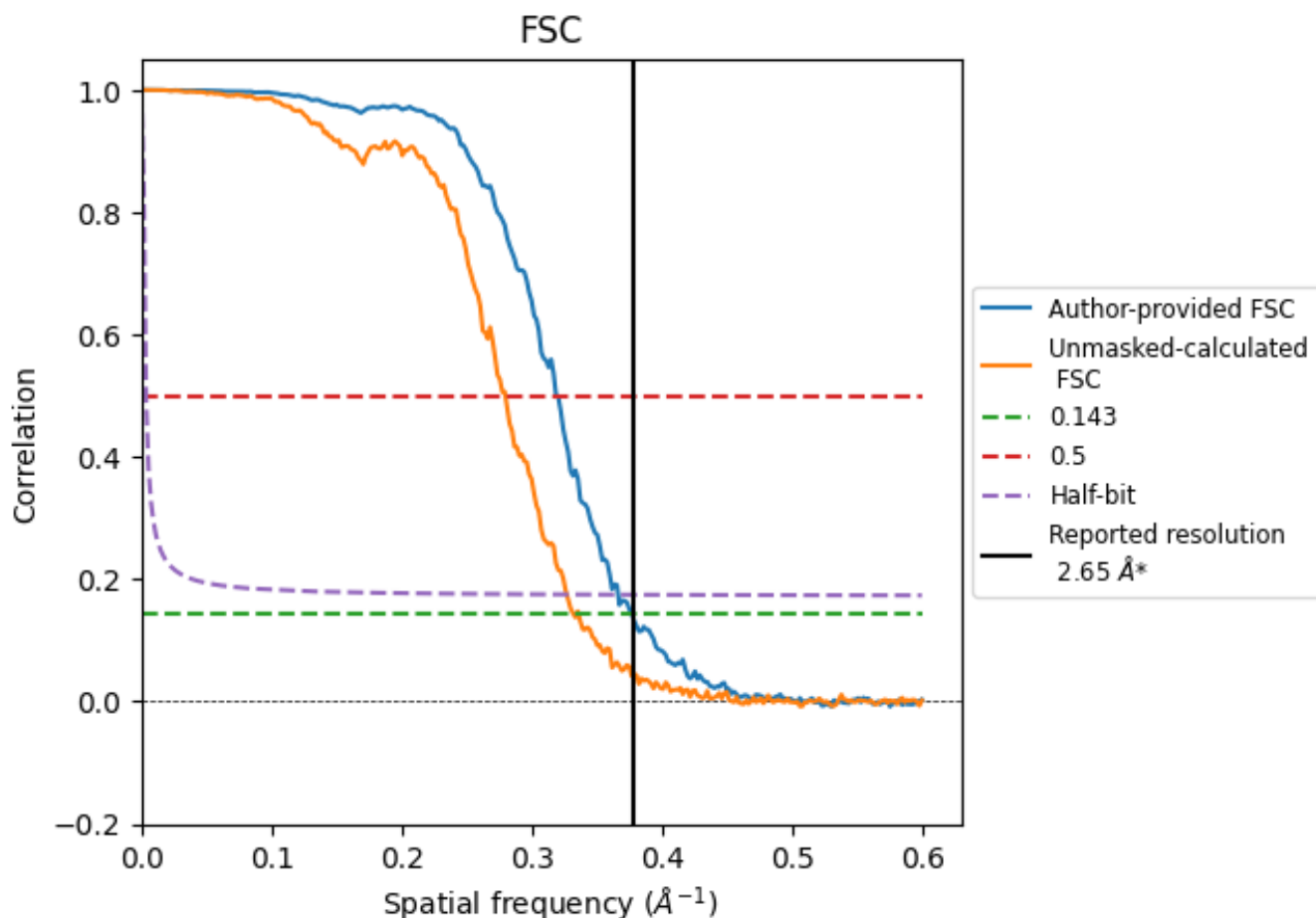


*Reported resolution corresponds to spatial frequency of 0.377 \AA^{-1}

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 0.377 \AA^{-1}

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.65	-	-
Author-provided FSC curve	2.65	3.13	2.73
Unmasked-calculated*	3.01	3.58	3.06

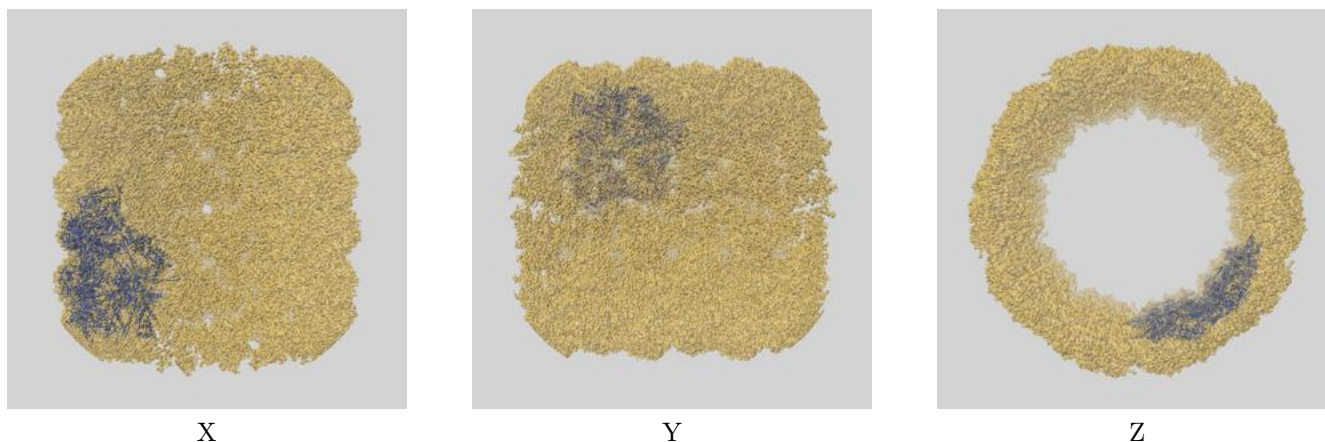
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.01 differs from the reported value 2.65 by more than 10 %

9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-17672 and PDB model 8PHR. Per-residue inclusion information can be found in section 3 on page 8.

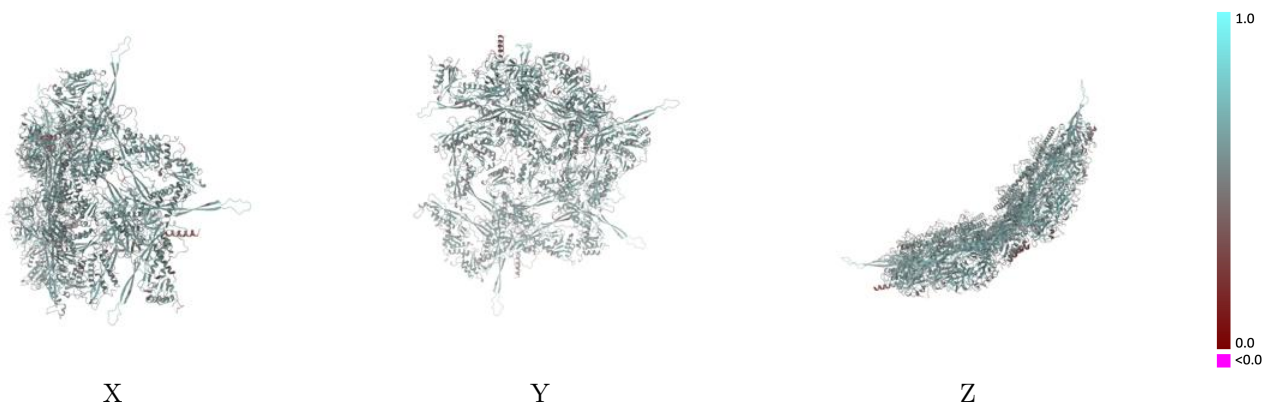
9.1 Map-model overlays

9.1.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 0.005 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](#)

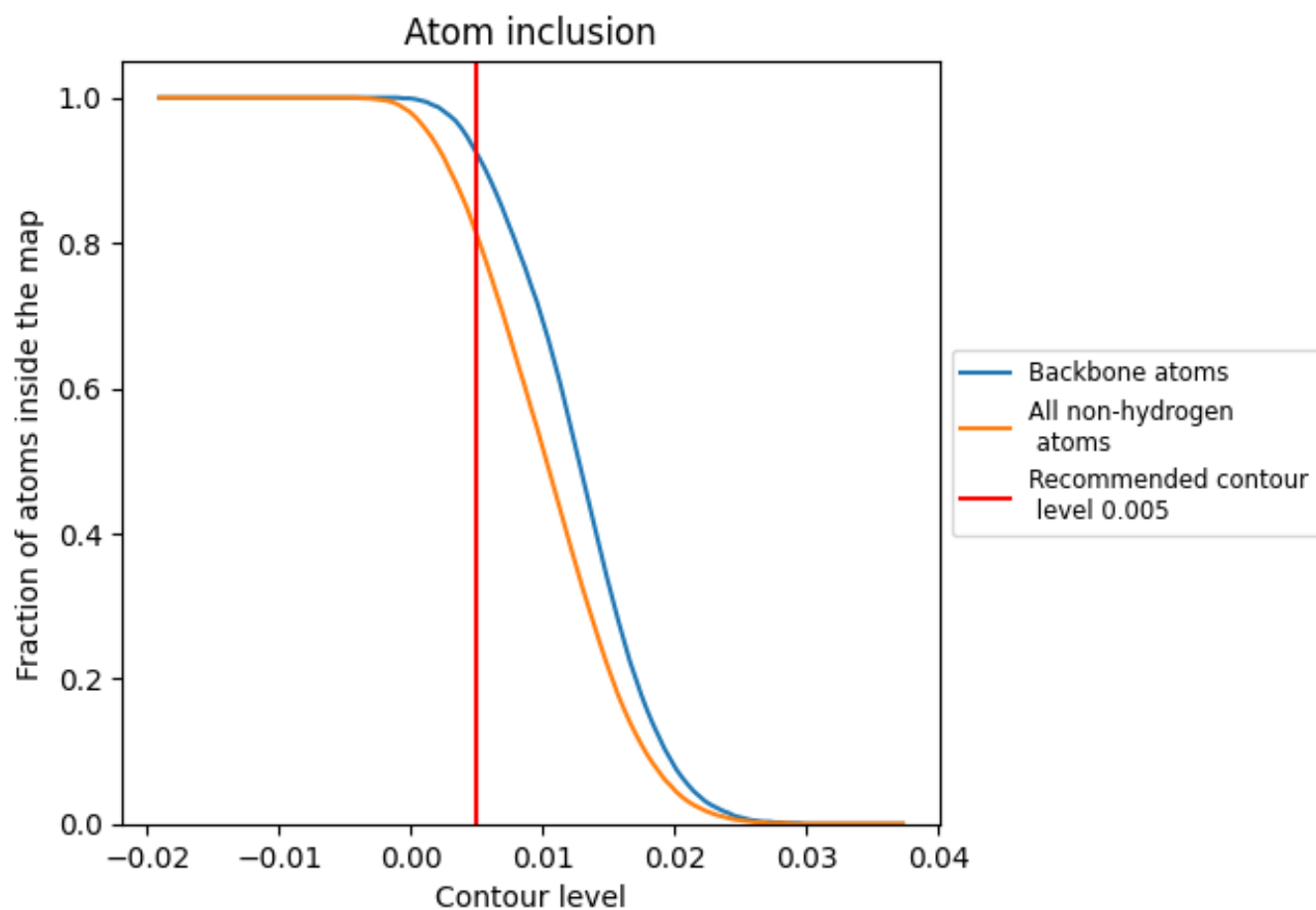


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 81% 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 (0.005) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8110	 0.5480
A	 0.8450	 0.5690
B	 0.8410	 0.5730
C	 0.8410	 0.5740
D	 0.8030	 0.5510
E	 0.8150	 0.5750
F	 0.7280	 0.5160
G	 0.7820	 0.5090
H	 0.7880	 0.5160
I	 0.7720	 0.5130
J	 0.8450	 0.5680
K	 0.8520	 0.5690
L	 0.8310	 0.5600
M	 0.7200	 0.5070
N	 0.8030	 0.5460
O	 0.8140	 0.5640
P	 0.8030	 0.5280
Q	 0.7950	 0.5170
R	 0.7950	 0.5210
S	 0.8370	 0.5610
T	 0.8290	 0.5630
U	 0.8180	 0.5530
V	 0.7780	 0.5360
W	 0.7870	 0.5600
X	 0.7080	 0.4990
Y	 0.7820	 0.4970
Z	 0.7620	 0.4900
a	 0.7910	 0.5080
b	 0.8360	 0.5540
c	 0.8460	 0.5640
d	 0.8080	 0.5440
e	 0.7030	 0.5010
f	 0.7930	 0.5330
g	 0.7870	 0.5490
h	 0.7670	 0.4750



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Chain	Atom inclusion	Q-score
i	 0.7690	 0.4960
j	 0.7650	 0.4970
k	 0.8180	 0.5470
l	 0.7260	 0.4870
m	 0.8110	 0.5470
n	 0.7550	 0.4870
o	 0.8340	 0.5510
p	 0.7330	 0.4920