



Full wwPDB EM Validation Report ⓘ

Mar 24, 2026 – 04:43 AM UTC

PDB ID : 9EUI / pdb_00009eui
EMDB ID : EMD-19971
Title : Cryo-EM structure of Staphylococcus aureus bacteriophage phi812 baseplate in the post-contraction state - complete
Authors : Binovsky, J.; Siborova, M.; Baska, R.; Pichel-Beleiro, A.; Skubnik, K.; Novacek, J.; van Raaij, M.J.; Plevka, P.
Deposited on : 2024-03-27
Resolution : 4.40 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

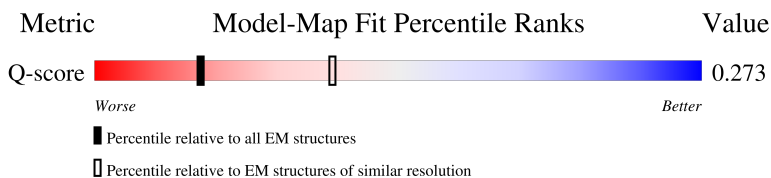
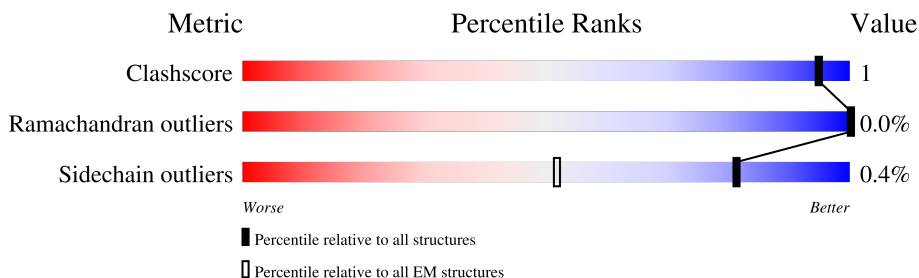
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.40 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.



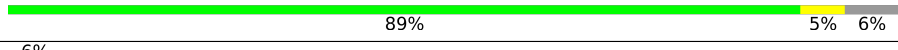
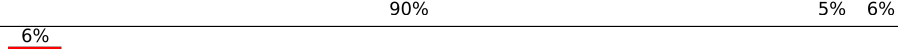
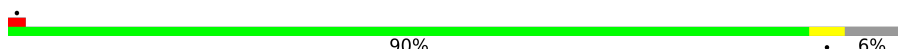
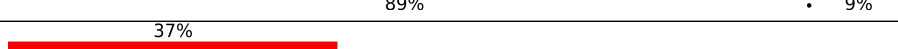
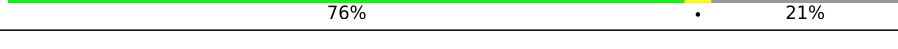
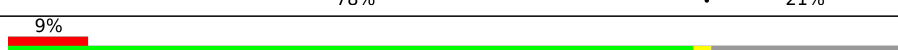
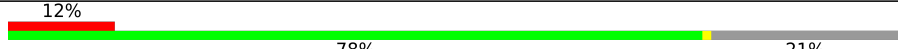
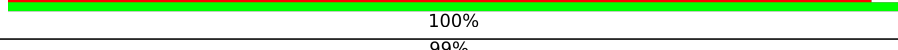
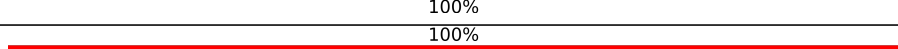
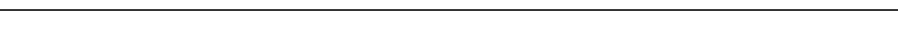
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	3132 (3.91 - 4.90)

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	234	95% .
2	B	348	97% .
2	C	348	95% 5%
3	D	1019	5% 36% . 63%

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Mol	Chain	Length	Quality of chain
4	E	587	
4	F	587	
4	G	587	
4	H	587	
4	I	587	
4	J	587	
5	K	194	
5	L	194	
5	M	194	
5	N	194	
5	O	194	
5	P	194	
6	Q	1152	
6	R	1152	
6	S	1152	
6	T	1152	
6	U	1152	
6	V	1152	
7	W	458	
7	X	458	
7	Y	458	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 92215 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 Baseplate wedge subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	234	Total	C	N	O	S	0	0
			1871	1174	314	377	6		

- Molecule 2 is a protein called Baseplate component.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	348	Total	C	N	O	S	0	0
			2760	1734	459	560	7		
2	C	347	Total	C	N	O	S	0	0
			2752	1729	458	559	6		

- Molecule 3 is a protein called TmpF.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	373	Total	C	N	O	S	0	0
			2252	1333	438	479	2		

- Molecule 4 is a protein called Major tail sheath protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	540	Total	C	N	O	S	0	0
			4213	2653	714	839	7		
4	F	553	Total	C	N	O	S	0	0
			4321	2723	734	857	7		
4	G	553	Total	C	N	O	S	0	0
			4321	2723	734	857	7		
4	H	553	Total	C	N	O	S	0	0
			4321	2723	734	857	7		
4	I	553	Total	C	N	O	S	0	0
			4321	2723	734	857	7		
4	J	553	Total	C	N	O	S	0	0
			4321	2723	734	857	7		

- Molecule 5 is a protein called Baseplate protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	K	176	Total	C	N	O	S	0	0
			1378	874	226	277	1		
5	L	176	Total	C	N	O	S	0	0
			1378	874	226	277	1		
5	M	176	Total	C	N	O	S	0	0
			1378	874	226	277	1		
5	N	176	Total	C	N	O	S	0	0
			1378	874	226	277	1		
5	O	176	Total	C	N	O	S	0	0
			1378	874	226	277	1		
5	P	176	Total	C	N	O	S	0	0
			1378	874	226	277	1		

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	174	ASP	-	expression tag	UNP A0A0U1WGQ5
K	175	PRO	-	expression tag	UNP A0A0U1WGQ5
K	176	ASN	-	expression tag	UNP A0A0U1WGQ5
K	177	SER	-	expression tag	UNP A0A0U1WGQ5
K	178	SER	-	expression tag	UNP A0A0U1WGQ5
K	179	SER	-	expression tag	UNP A0A0U1WGQ5
K	180	VAL	-	expression tag	UNP A0A0U1WGQ5
K	181	ASP	-	expression tag	UNP A0A0U1WGQ5
K	182	LYS	-	expression tag	UNP A0A0U1WGQ5
K	183	LEU	-	expression tag	UNP A0A0U1WGQ5
K	184	ALA	-	expression tag	UNP A0A0U1WGQ5
K	185	ALA	-	expression tag	UNP A0A0U1WGQ5
K	186	ALA	-	expression tag	UNP A0A0U1WGQ5
K	187	LEU	-	expression tag	UNP A0A0U1WGQ5
K	188	GLU	-	expression tag	UNP A0A0U1WGQ5
K	189	HIS	-	expression tag	UNP A0A0U1WGQ5
K	190	HIS	-	expression tag	UNP A0A0U1WGQ5
K	191	HIS	-	expression tag	UNP A0A0U1WGQ5
K	192	HIS	-	expression tag	UNP A0A0U1WGQ5
K	193	HIS	-	expression tag	UNP A0A0U1WGQ5
K	194	HIS	-	expression tag	UNP A0A0U1WGQ5
L	174	ASP	-	expression tag	UNP A0A0U1WGQ5
L	175	PRO	-	expression tag	UNP A0A0U1WGQ5
L	176	ASN	-	expression tag	UNP A0A0U1WGQ5
L	177	SER	-	expression tag	UNP A0A0U1WGQ5
L	178	SER	-	expression tag	UNP A0A0U1WGQ5

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Chain	Residue	Modelled	Actual	Comment	Reference
L	179	SER	-	expression tag	UNP A0A0U1WGGQ5
L	180	VAL	-	expression tag	UNP A0A0U1WGGQ5
L	181	ASP	-	expression tag	UNP A0A0U1WGGQ5
L	182	LYS	-	expression tag	UNP A0A0U1WGGQ5
L	183	LEU	-	expression tag	UNP A0A0U1WGGQ5
L	184	ALA	-	expression tag	UNP A0A0U1WGGQ5
L	185	ALA	-	expression tag	UNP A0A0U1WGGQ5
L	186	ALA	-	expression tag	UNP A0A0U1WGGQ5
L	187	LEU	-	expression tag	UNP A0A0U1WGGQ5
L	188	GLU	-	expression tag	UNP A0A0U1WGGQ5
L	189	HIS	-	expression tag	UNP A0A0U1WGGQ5
L	190	HIS	-	expression tag	UNP A0A0U1WGGQ5
L	191	HIS	-	expression tag	UNP A0A0U1WGGQ5
L	192	HIS	-	expression tag	UNP A0A0U1WGGQ5
L	193	HIS	-	expression tag	UNP A0A0U1WGGQ5
L	194	HIS	-	expression tag	UNP A0A0U1WGGQ5
M	174	ASP	-	expression tag	UNP A0A0U1WGGQ5
M	175	PRO	-	expression tag	UNP A0A0U1WGGQ5
M	176	ASN	-	expression tag	UNP A0A0U1WGGQ5
M	177	SER	-	expression tag	UNP A0A0U1WGGQ5
M	178	SER	-	expression tag	UNP A0A0U1WGGQ5
M	179	SER	-	expression tag	UNP A0A0U1WGGQ5
M	180	VAL	-	expression tag	UNP A0A0U1WGGQ5
M	181	ASP	-	expression tag	UNP A0A0U1WGGQ5
M	182	LYS	-	expression tag	UNP A0A0U1WGGQ5
M	183	LEU	-	expression tag	UNP A0A0U1WGGQ5
M	184	ALA	-	expression tag	UNP A0A0U1WGGQ5
M	185	ALA	-	expression tag	UNP A0A0U1WGGQ5
M	186	ALA	-	expression tag	UNP A0A0U1WGGQ5
M	187	LEU	-	expression tag	UNP A0A0U1WGGQ5
M	188	GLU	-	expression tag	UNP A0A0U1WGGQ5
M	189	HIS	-	expression tag	UNP A0A0U1WGGQ5
M	190	HIS	-	expression tag	UNP A0A0U1WGGQ5
M	191	HIS	-	expression tag	UNP A0A0U1WGGQ5
M	192	HIS	-	expression tag	UNP A0A0U1WGGQ5
M	193	HIS	-	expression tag	UNP A0A0U1WGGQ5
M	194	HIS	-	expression tag	UNP A0A0U1WGGQ5
N	174	ASP	-	expression tag	UNP A0A0U1WGGQ5
N	175	PRO	-	expression tag	UNP A0A0U1WGGQ5
N	176	ASN	-	expression tag	UNP A0A0U1WGGQ5
N	177	SER	-	expression tag	UNP A0A0U1WGGQ5
N	178	SER	-	expression tag	UNP A0A0U1WGGQ5

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Chain	Residue	Modelled	Actual	Comment	Reference
N	179	SER	-	expression tag	UNP A0A0U1WGGQ5
N	180	VAL	-	expression tag	UNP A0A0U1WGGQ5
N	181	ASP	-	expression tag	UNP A0A0U1WGGQ5
N	182	LYS	-	expression tag	UNP A0A0U1WGGQ5
N	183	LEU	-	expression tag	UNP A0A0U1WGGQ5
N	184	ALA	-	expression tag	UNP A0A0U1WGGQ5
N	185	ALA	-	expression tag	UNP A0A0U1WGGQ5
N	186	ALA	-	expression tag	UNP A0A0U1WGGQ5
N	187	LEU	-	expression tag	UNP A0A0U1WGGQ5
N	188	GLU	-	expression tag	UNP A0A0U1WGGQ5
N	189	HIS	-	expression tag	UNP A0A0U1WGGQ5
N	190	HIS	-	expression tag	UNP A0A0U1WGGQ5
N	191	HIS	-	expression tag	UNP A0A0U1WGGQ5
N	192	HIS	-	expression tag	UNP A0A0U1WGGQ5
N	193	HIS	-	expression tag	UNP A0A0U1WGGQ5
N	194	HIS	-	expression tag	UNP A0A0U1WGGQ5
O	174	ASP	-	expression tag	UNP A0A0U1WGGQ5
O	175	PRO	-	expression tag	UNP A0A0U1WGGQ5
O	176	ASN	-	expression tag	UNP A0A0U1WGGQ5
O	177	SER	-	expression tag	UNP A0A0U1WGGQ5
O	178	SER	-	expression tag	UNP A0A0U1WGGQ5
O	179	SER	-	expression tag	UNP A0A0U1WGGQ5
O	180	VAL	-	expression tag	UNP A0A0U1WGGQ5
O	181	ASP	-	expression tag	UNP A0A0U1WGGQ5
O	182	LYS	-	expression tag	UNP A0A0U1WGGQ5
O	183	LEU	-	expression tag	UNP A0A0U1WGGQ5
O	184	ALA	-	expression tag	UNP A0A0U1WGGQ5
O	185	ALA	-	expression tag	UNP A0A0U1WGGQ5
O	186	ALA	-	expression tag	UNP A0A0U1WGGQ5
O	187	LEU	-	expression tag	UNP A0A0U1WGGQ5
O	188	GLU	-	expression tag	UNP A0A0U1WGGQ5
O	189	HIS	-	expression tag	UNP A0A0U1WGGQ5
O	190	HIS	-	expression tag	UNP A0A0U1WGGQ5
O	191	HIS	-	expression tag	UNP A0A0U1WGGQ5
O	192	HIS	-	expression tag	UNP A0A0U1WGGQ5
O	193	HIS	-	expression tag	UNP A0A0U1WGGQ5
O	194	HIS	-	expression tag	UNP A0A0U1WGGQ5
P	174	ASP	-	expression tag	UNP A0A0U1WGGQ5
P	175	PRO	-	expression tag	UNP A0A0U1WGGQ5
P	176	ASN	-	expression tag	UNP A0A0U1WGGQ5
P	177	SER	-	expression tag	UNP A0A0U1WGGQ5
P	178	SER	-	expression tag	UNP A0A0U1WGGQ5

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Chain	Residue	Modelled	Actual	Comment	Reference
P	179	SER	-	expression tag	UNP A0A0U1WGQ5
P	180	VAL	-	expression tag	UNP A0A0U1WGQ5
P	181	ASP	-	expression tag	UNP A0A0U1WGQ5
P	182	LYS	-	expression tag	UNP A0A0U1WGQ5
P	183	LEU	-	expression tag	UNP A0A0U1WGQ5
P	184	ALA	-	expression tag	UNP A0A0U1WGQ5
P	185	ALA	-	expression tag	UNP A0A0U1WGQ5
P	186	ALA	-	expression tag	UNP A0A0U1WGQ5
P	187	LEU	-	expression tag	UNP A0A0U1WGQ5
P	188	GLU	-	expression tag	UNP A0A0U1WGQ5
P	189	HIS	-	expression tag	UNP A0A0U1WGQ5
P	190	HIS	-	expression tag	UNP A0A0U1WGQ5
P	191	HIS	-	expression tag	UNP A0A0U1WGQ5
P	192	HIS	-	expression tag	UNP A0A0U1WGQ5
P	193	HIS	-	expression tag	UNP A0A0U1WGQ5
P	194	HIS	-	expression tag	UNP A0A0U1WGQ5

- Molecule 6 is a protein called DUF4815 domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	Q	908	Total	C	N	O	S	0	0
			7166	4511	1184	1453	18		
6	R	908	Total	C	N	O	S	0	0
			7166	4511	1184	1453	18		
6	S	908	Total	C	N	O	S	0	0
			7166	4511	1184	1453	18		
6	T	908	Total	C	N	O	S	0	0
			7166	4511	1184	1453	18		
6	U	908	Total	C	N	O	S	0	0
			7165	4510	1184	1453	18		
6	V	908	Total	C	N	O	S	0	0
			7166	4511	1184	1453	18		

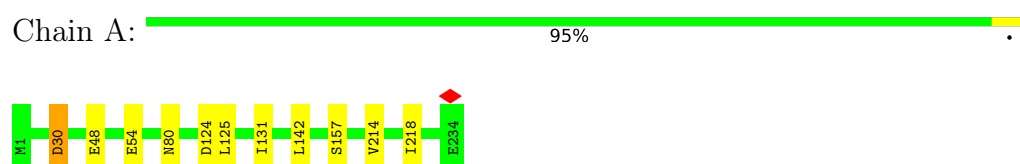
- Molecule 7 is a protein called Receptor binding protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	W	458	Total	C	N	O	0	0
			1833	916	458	459		
7	X	458	Total	C	N	O	0	0
			1833	916	458	459		
7	Y	458	Total	C	N	O	0	0
			1833	916	458	459		

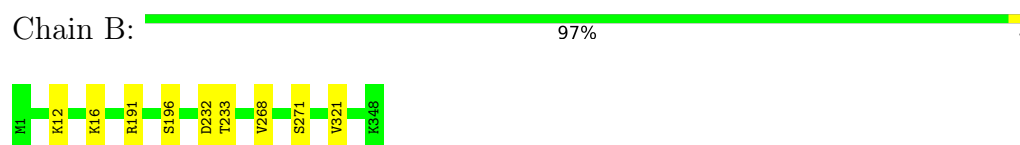
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

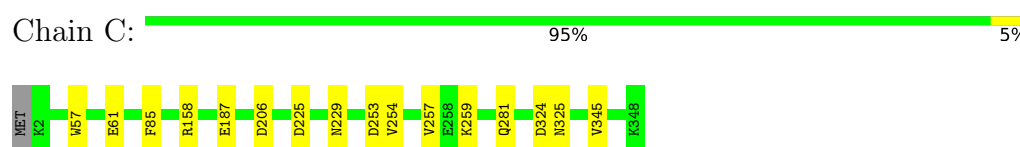
- Molecule 1: Baseplate wedge subunit



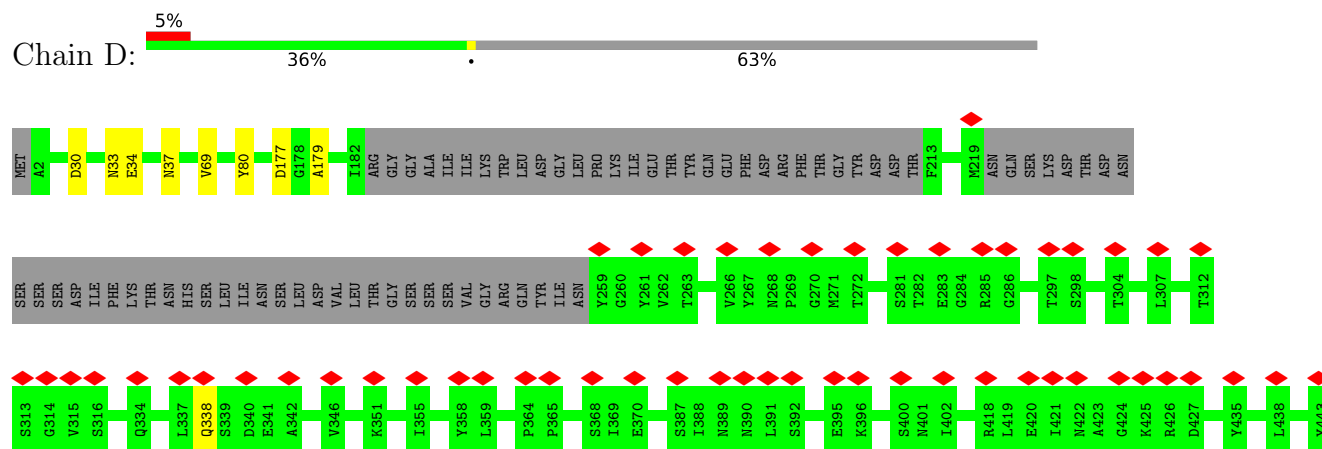
- Molecule 2: Baseplate component

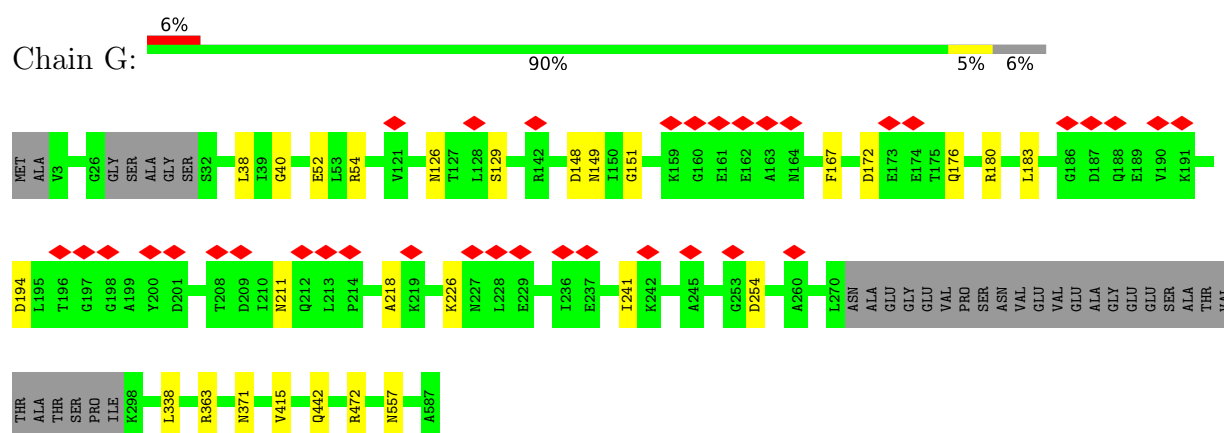


- Molecule 2: Baseplate component

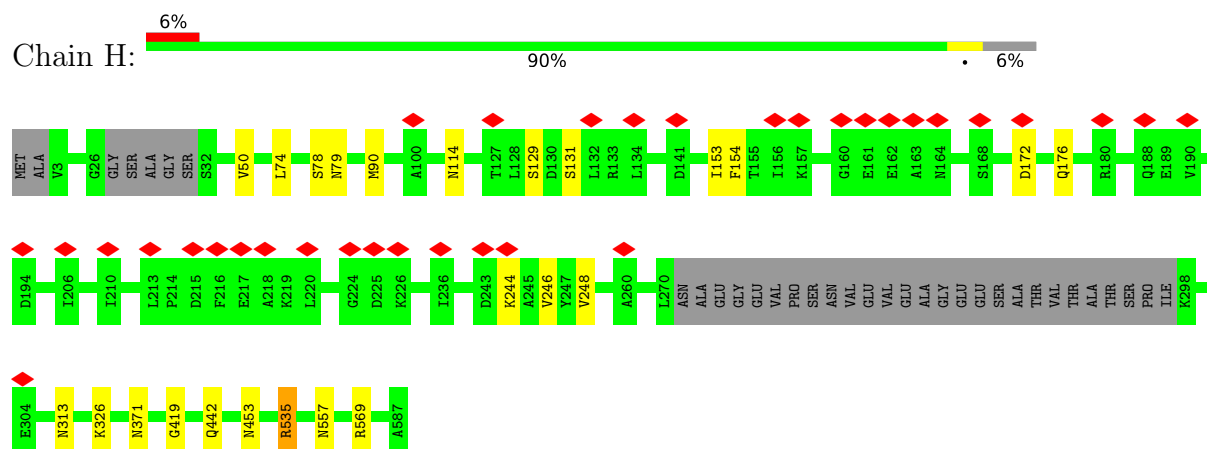


- Molecule 3: TmpF

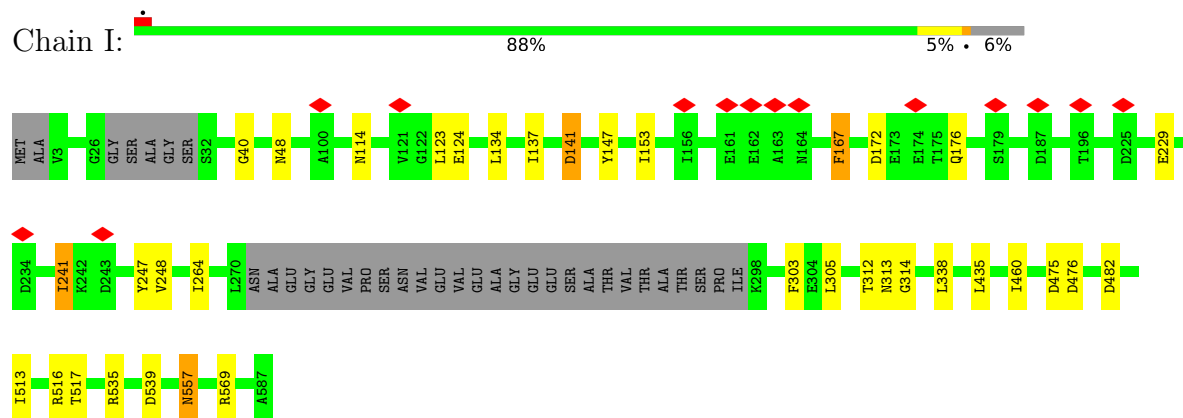




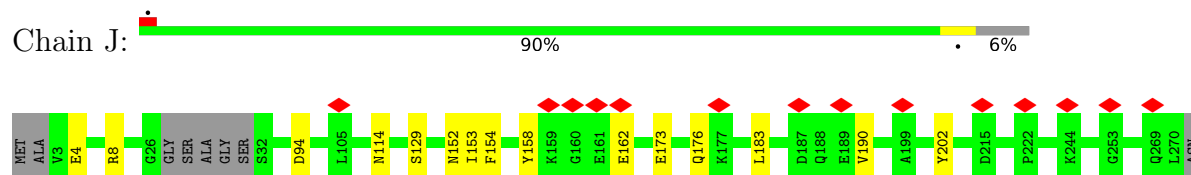
- Molecule 4: Major tail sheath protein

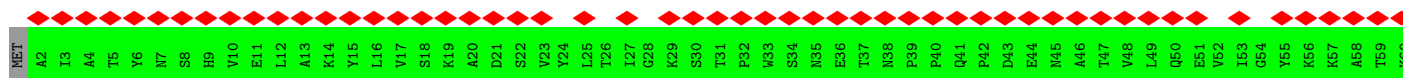


- Molecule 4: Major tail sheath protein



- Molecule 4: Major tail sheath protein





21%

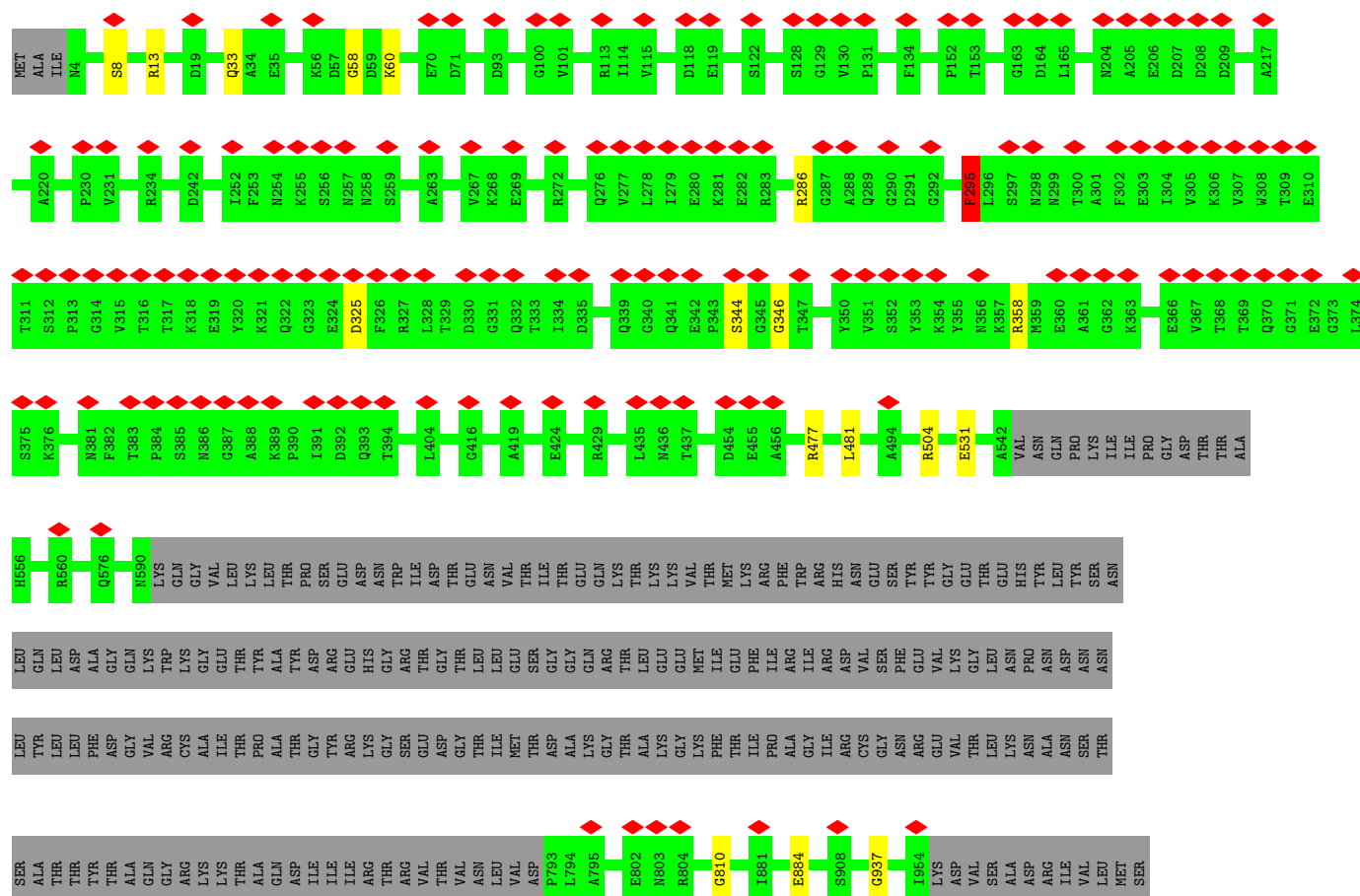
21%

21%

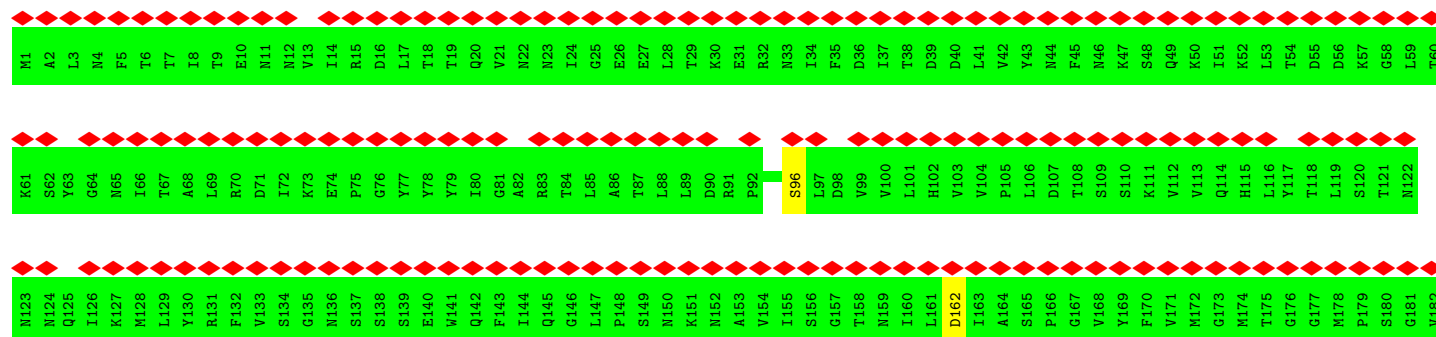


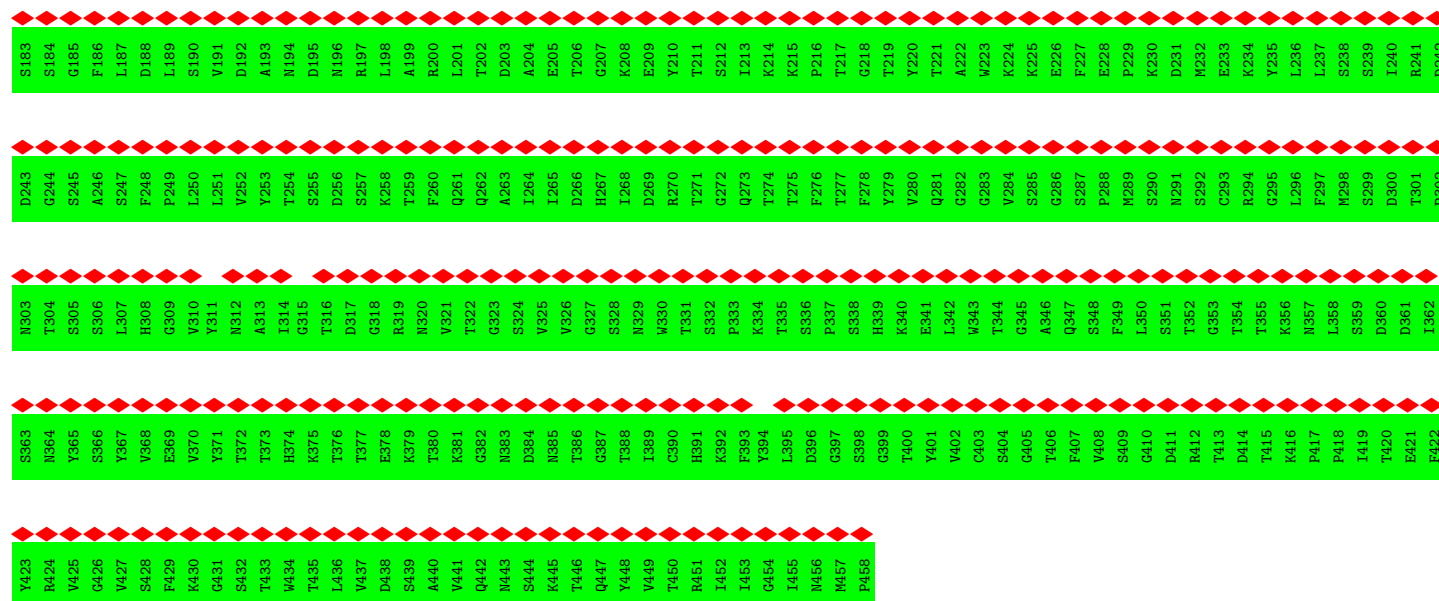


Chain V: 14% 77% 21%



Chain W:





- Molecule 7: Receptor binding protein



Chain Y:



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C6	Depositor
Number of particles used	25203	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 ($e^-/\text{\AA}^2$)	42	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	130000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.052	Depositor
Minimum map value	-0.023	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.006	Depositor
Map size (Å)	761.04004, 761.04004, 761.04004	wwPDB
Map dimensions	720, 720, 720	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.057, 1.057, 1.057	Depositor

5 Model quality

5.1 Standard geometry

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	0.23	0/1902	0.49	0/2572
2	B	0.41	0/2803	0.74	0/3794
2	C	0.21	0/2795	0.44	0/3784
3	D	0.47	0/2281	0.87	0/3004
4	E	0.24	0/4280	0.48	0/5777
4	F	0.33	0/4392	0.63	0/5930
4	G	0.35	0/4392	0.64	1/5930 (0.0%)
4	H	0.32	0/4392	0.62	1/5930 (0.0%)
4	I	0.28	0/4392	0.56	0/5930
4	J	0.21	0/4392	0.45	0/5930
5	K	0.70	0/1407	1.21	3/1914 (0.2%)
5	L	0.73	0/1407	1.31	4/1914 (0.2%)
5	M	0.80	0/1407	1.31	2/1914 (0.1%)
5	N	0.81	0/1407	1.31	1/1914 (0.1%)
5	O	0.86	0/1407	1.35	2/1914 (0.1%)
5	P	0.84	0/1407	1.32	0/1914
6	Q	0.28	0/7308	0.52	1/9911 (0.0%)
6	R	0.29	0/7308	0.54	1/9911 (0.0%)
6	S	0.45	0/7308	0.77	0/9911
6	T	0.54	0/7308	0.91	6/9911 (0.1%)
6	U	0.50	0/7305	0.82	2/9905 (0.0%)
6	V	0.58	0/7308	0.95	4/9911 (0.0%)
7	W	0.85	0/1832	1.28	0/2287
7	X	0.86	0/1832	1.27	0/2287
7	Y	0.85	0/1832	1.28	0/2287
All	All	0.48	0/93804	0.81	28/126386 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

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Mol	Chain	#Chirality outliers	#Planarity outliers
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Mol	Chain	#Chirality outliers	#Planarity outliers
4	G	0	1
4	H	0	1
5	O	0	1
6	S	0	2
6	T	0	2
6	V	0	2
All	All	0	9

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	V	295	PHE	CA-CB-CG	-8.65	105.15	113.80
5	M	169	PHE	CA-CB-CG	7.13	120.94	113.80
5	O	169	PHE	CA-CB-CG	7.04	120.84	113.80
5	L	169	PHE	CA-CB-CG	6.91	120.71	113.80
6	U	438	ASP	CA-CB-CG	6.81	119.41	112.60
6	T	1144	ARG	NE-CZ-NH2	6.78	125.30	119.20
5	L	38	ASN	CA-C-N	6.77	124.60	119.66
5	L	38	ASN	C-N-CA	6.77	124.60	119.66
6	T	477	ARG	NE-CZ-NH2	6.63	125.17	119.20
4	H	535	ARG	NE-CZ-NH2	6.42	124.98	119.20
5	N	169	PHE	CA-CB-CG	6.30	120.10	113.80
5	M	65	ARG	NE-CZ-NH2	6.22	124.80	119.20
5	L	176	ASN	CA-CB-CG	6.18	118.78	112.60
6	V	504	ARG	NE-CZ-NH2	6.11	124.70	119.20
6	T	286	ARG	NE-CZ-NH2	6.10	124.69	119.20
6	Q	1144	ARG	NE-CZ-NH2	5.78	124.40	119.20
6	U	477	ARG	NE-CZ-NH2	5.77	124.39	119.20
5	O	65	ARG	NE-CZ-NH2	5.61	124.25	119.20
6	T	272	ARG	NE-CZ-NH2	5.61	124.24	119.20
4	G	472	ARG	NE-CZ-NH2	5.59	124.23	119.20
5	K	135	PHE	CA-CB-CG	5.55	119.35	113.80
5	K	74	ASN	CA-CB-CG	5.48	118.08	112.60
6	V	477	ARG	NE-CZ-NH2	5.36	124.02	119.20
6	T	504	ARG	NE-CZ-NH2	5.21	123.89	119.20
6	R	1128	ARG	NE-CZ-NH2	5.21	123.89	119.20
6	T	57	ASP	CA-CB-CG	5.06	117.66	112.60
5	K	169	PHE	CA-CB-CG	5.05	118.86	113.80
6	V	1086	ASP	CA-CB-CG	5.03	117.63	112.60

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	G	363	ARG	Sidechain
4	H	535	ARG	Sidechain
5	O	158	ASN	Peptide
6	S	272	ARG	Sidechain
6	S	286	ARG	Sidechain
6	T	286	ARG	Sidechain
6	T	327	ARG	Sidechain
6	V	286	ARG	Sidechain
6	V	295	PHE	Sidechain

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	1871	0	1828	9	0
2	B	2760	0	2729	6	0
2	C	2752	0	2717	8	0
3	D	2252	0	1652	13	0
4	E	4213	0	4167	21	0
4	F	4321	0	4280	17	0
4	G	4321	0	4280	16	0
4	H	4321	0	4280	14	0
4	I	4321	0	4280	22	0
4	J	4321	0	4280	14	0
5	K	1378	0	1361	7	0
5	L	1378	0	1361	0	0
5	M	1378	0	1361	0	0
5	N	1378	0	1361	1	0
5	O	1378	0	1361	1	0
5	P	1378	0	1361	0	0
6	Q	7166	0	6971	14	0
6	R	7166	0	6971	17	0
6	S	7166	0	6971	7	0
6	T	7166	0	6971	10	0
6	U	7165	0	6969	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	V	7166	0	6971	20	0
7	W	1833	0	520	13	0
7	X	1833	0	520	0	0
7	Y	1833	0	520	0	0
All	All	92215	0	86043	189	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 1.

All (189) 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)
6:V:295:PHE:CE2	7:W:96:SER:N	1.85	1.39
6:V:344:SER:CB	7:W:162:ASP:CA	2.31	1.09
6:V:295:PHE:CD2	7:W:96:SER:N	2.13	1.06
6:V:344:SER:OG	7:W:162:ASP:CA	2.11	0.98
3:D:177:ASP:CG	5:K:168:ARG:HH11	1.74	0.94
6:V:295:PHE:CD2	7:W:96:SER:CA	2.50	0.94
3:D:177:ASP:OD1	5:K:168:ARG:NH1	2.02	0.91
6:V:344:SER:HB3	7:W:162:ASP:CA	2.00	0.90
6:V:344:SER:HB3	7:W:162:ASP:C	2.00	0.86
6:V:295:PHE:HE2	7:W:96:SER:H	0.90	0.84
6:U:1077:THR:CA	6:U:1078:LYS:N	2.41	0.83
6:V:344:SER:HB3	7:W:162:ASP:O	1.83	0.79
6:V:295:PHE:CE2	7:W:96:SER:CA	2.68	0.76
3:D:33:ASN:OD1	3:D:37:ASN:ND2	2.20	0.74
2:B:191:ARG:NH1	2:C:187:GLU:OE1	2.21	0.74
3:D:179:ALA:HB1	5:K:80:TYR:HA	1.71	0.71
6:U:1077:THR:O	6:U:1078:LYS:N	2.26	0.69
3:D:338:GLN:O	6:T:4:ASN:HB3	1.93	0.69
6:U:1077:THR:CA	6:U:1077:THR:O	2.41	0.68
2:C:85:PHE:O	2:C:158:ARG:NH2	2.26	0.68
4:G:557:ASN:OD1	4:H:569:ARG:NH2	2.25	0.68
6:U:83:ILE:O	6:U:88:ARG:NH2	2.26	0.68
6:S:83:ILE:O	6:S:88:ARG:NH2	2.26	0.68
6:Q:83:ILE:O	6:Q:88:ARG:NH2	2.27	0.67
4:F:371:ASN:OD1	4:F:442:GLN:NE2	2.29	0.65
3:D:177:ASP:CG	5:K:168:ARG:NH1	2.51	0.65
4:E:557:ASN:OD1	4:F:569:ARG:NH2	2.31	0.64
4:I:141:ASP:OD1	4:I:141:ASP:N	2.27	0.64
6:Q:88:ARG:NH1	6:Q:143:GLU:OE1	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:Q:504:ARG:NH1	6:S:480:ASN:OD1	2.32	0.63
6:Q:480:ASN:OD1	6:R:504:ARG:NH1	2.32	0.63
4:H:371:ASN:OD1	4:H:442:GLN:NE2	2.33	0.62
4:I:124:GLU:N	4:I:124:GLU:OE1	2.33	0.61
6:V:295:PHE:HE2	7:W:96:SER:N	1.53	0.60
4:J:329:HIS:ND1	4:J:360:GLU:OE2	2.35	0.60
4:I:516:ARG:NH1	4:J:428:GLU:OE2	2.35	0.59
4:J:129:SER:O	4:J:152:ASN:ND2	2.36	0.59
2:C:57:TRP:NE1	2:C:61:GLU:OE1	2.35	0.59
4:H:74:LEU:O	4:H:78:SER:N	2.36	0.58
4:E:386:ASN:O	4:E:479:THR:OG1	2.22	0.58
4:E:70:ASP:OD1	4:E:404:ARG:NH1	2.36	0.58
2:B:12:LYS:NZ	3:D:34:GLU:OE2	2.36	0.58
1:A:80:ASN:ND2	4:E:83:THR:OG1	2.37	0.57
6:V:344:SER:CB	7:W:162:ASP:O	2.52	0.57
6:R:480:ASN:OD1	6:S:504:ARG:NH1	2.37	0.57
4:J:4:GLU:OE2	4:J:8:ARG:NH2	2.38	0.57
3:D:338:GLN:O	6:T:4:ASN:CB	2.53	0.57
4:J:94:ASP:OD1	4:J:323:LYS:NZ	2.38	0.56
6:V:344:SER:HG	7:W:162:ASP:CA	2.18	0.55
4:J:158:TYR:OH	4:J:162:GLU:N	2.38	0.55
1:A:131:ILE:HG21	4:E:22:THR:HG23	1.89	0.55
3:D:179:ALA:CB	5:K:80:TYR:HA	2.37	0.54
4:F:482:ASP:N	4:F:482:ASP:OD1	2.39	0.54
4:F:172:ASP:O	4:F:176:GLN:N	2.40	0.54
4:J:153:ILE:HG23	4:J:154:PHE:HD1	1.74	0.53
6:R:216:ASP:OD1	6:R:217:ALA:N	2.41	0.53
4:G:371:ASN:OD1	4:G:442:GLN:NE2	2.41	0.53
4:I:535:ARG:NH1	4:I:539:ASP:OD2	2.41	0.53
4:I:229:GLU:N	4:I:229:GLU:OE1	2.41	0.53
4:H:153:ILE:HG23	4:H:154:PHE:HD1	1.72	0.53
4:E:172:ASP:O	4:E:176:GLN:N	2.40	0.53
4:E:148:ASP:OD1	4:E:149:ASN:N	2.41	0.52
4:G:148:ASP:OD1	4:G:149:ASN:N	2.42	0.52
4:H:557:ASN:OD1	4:I:569:ARG:NH2	2.42	0.52
4:H:153:ILE:HG22	4:H:248:VAL:O	2.09	0.52
4:E:164:ASN:OD1	4:E:165:ALA:N	2.42	0.52
4:E:517:THR:N	4:F:569:ARG:O	2.41	0.52
4:E:34:LYS:NZ	4:E:495:GLU:OE2	2.31	0.51
4:E:206:ILE:HG22	4:E:210:ILE:CD1	2.39	0.51
4:E:153:ILE:HG23	4:E:154:PHE:CD1	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:167:PHE:HD1	4:G:241:ILE:HD11	1.75	0.51
4:E:464:ARG:O	4:E:466:ARG:NH2	2.43	0.51
6:Q:477:ARG:NH1	6:R:479:ASP:OD1	2.40	0.50
2:B:232:ASP:OD1	2:B:233:THR:N	2.44	0.50
6:R:239:LYS:NZ	6:R:454:ASP:OD1	2.36	0.50
6:R:818:ASP:OD1	6:R:820:GLN:N	2.43	0.50
4:H:172:ASP:O	4:H:176:GLN:N	2.42	0.50
6:Q:426:ASN:OD1	6:Q:427:ILE:N	2.45	0.50
4:I:513:ILE:HG23	4:J:572:LYS:HA	1.92	0.50
1:A:30:ASP:N	1:A:30:ASP:OD1	2.45	0.49
4:F:123:LEU:HD12	4:F:303:PHE:CE2	2.47	0.49
4:H:50:VAL:HG22	4:H:90:MET:HE3	1.95	0.49
1:A:124:ASP:OD1	1:A:125:LEU:N	2.46	0.49
2:B:16:LYS:NZ	3:D:30:ASP:OD2	2.45	0.49
1:A:131:ILE:HG21	4:E:22:THR:CG2	2.44	0.48
2:B:191:ARG:O	2:B:196:SER:OG	2.30	0.48
6:V:810:GLY:N	6:V:937:GLY:O	2.45	0.48
4:E:509:GLU:HA	4:E:513:ILE:HD12	1.95	0.48
4:F:148:ASP:OD1	4:F:149:ASN:N	2.46	0.48
4:F:124:GLU:O	4:F:133:ARG:N	2.46	0.48
4:H:90:MET:HE1	4:H:326:LYS:HB2	1.96	0.48
6:Q:479:ASP:OD1	6:S:477:ARG:NH1	2.45	0.48
4:E:206:ILE:HG22	4:E:210:ILE:HD11	1.95	0.48
4:H:114:ASN:N	4:H:313:ASN:OD1	2.47	0.48
4:I:557:ASN:N	4:I:557:ASN:OD1	2.47	0.48
4:I:123:LEU:HD12	4:I:303:PHE:CE2	2.49	0.47
6:Q:239:LYS:NZ	6:Q:454:ASP:OD1	2.42	0.47
1:A:142:LEU:O	1:A:157:SER:OG	2.31	0.47
3:D:177:ASP:OD2	5:K:168:ARG:NH1	2.48	0.47
4:I:517:THR:N	4:J:569:ARG:O	2.45	0.47
4:G:167:PHE:CB	4:G:183:LEU:HD23	2.45	0.47
4:I:40:GLY:HA3	4:I:338:LEU:HD12	1.97	0.47
1:A:54:GLU:OE1	1:A:54:GLU:N	2.42	0.46
6:T:818:ASP:OD1	6:T:820:GLN:N	2.43	0.46
4:G:151:GLY:O	4:G:226:LYS:NZ	2.36	0.46
4:I:435:LEU:HD22	4:I:460:ILE:HD13	1.96	0.46
4:I:475:ASP:OD1	4:I:476:ASP:N	2.45	0.46
4:G:211:ASN:ND2	4:G:218:ALA:O	2.48	0.46
4:I:312:THR:HG22	4:I:314:GLY:H	1.81	0.46
4:I:482:ASP:OD1	4:I:482:ASP:N	2.49	0.46
6:R:20:ARG:NH1	6:S:120:ASP:OD2	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:R:88:ARG:NH1	6:R:143:GLU:OE1	2.49	0.46
4:E:516:ARG:NH1	4:F:426:ILE:O	2.49	0.46
2:C:225:ASP:OD1	2:C:229:ASN:N	2.49	0.46
4:I:172:ASP:O	4:I:176:GLN:N	2.42	0.46
4:H:153:ILE:HG23	4:H:154:PHE:CD1	2.50	0.45
4:G:126:ASN:ND2	4:G:129:SER:OG	2.46	0.45
6:T:20:ARG:NH1	6:U:120:ASP:OD2	2.44	0.45
4:F:375:GLU:OE1	4:F:375:GLU:N	2.47	0.45
2:C:324:ASP:OD1	2:C:325:ASN:N	2.49	0.45
6:Q:1084:SER:OG	6:Q:1086:ASP:OD1	2.34	0.45
4:G:40:GLY:HA3	4:G:338:LEU:HD12	1.99	0.45
4:G:167:PHE:CD1	4:G:241:ILE:HD11	2.52	0.45
3:D:179:ALA:HB1	5:K:81:GLY:H	1.82	0.45
4:H:129:SER:OG	4:H:131:SER:OG	2.33	0.45
3:D:69:VAL:O	3:D:80:TYR:OH	2.34	0.45
6:Q:810:GLY:N	6:Q:937:GLY:O	2.45	0.44
4:E:74:LEU:O	4:E:78:SER:N	2.50	0.44
4:G:180:ARG:NE	4:G:194:ASP:OD1	2.42	0.44
4:J:114:ASN:OD1	4:J:313:ASN:N	2.49	0.44
4:I:153:ILE:HG22	4:I:248:VAL:O	2.18	0.44
4:I:114:ASN:OD1	4:I:313:ASN:N	2.51	0.44
4:I:312:THR:HG22	4:I:314:GLY:N	2.32	0.44
4:J:183:LEU:O	4:J:190:VAL:N	2.48	0.44
6:R:92:ASN:ND2	6:R:145:MET:O	2.51	0.44
4:E:577:SER:C	4:E:578:LEU:HD12	2.43	0.44
6:R:426:ASN:OD1	6:R:427:ILE:N	2.49	0.44
4:I:134:LEU:N	4:I:147:TYR:O	2.51	0.43
4:H:79:ASN:ND2	4:H:419:GLY:O	2.50	0.43
6:R:518:ASP:OD1	6:T:347:THR:HB	2.18	0.43
6:T:92:ASN:ND2	6:T:145:MET:O	2.51	0.43
6:Q:50:LEU:HA	6:R:87:ILE:HD12	2.01	0.43
1:A:214:VAL:HG12	1:A:218:ILE:HD12	2.00	0.43
4:I:137:ILE:HD12	4:I:305:LEU:HD21	2.01	0.43
6:T:122:SER:O	6:V:13:ARG:NH2	2.51	0.43
4:F:153:ILE:HG22	4:F:248:VAL:O	2.19	0.43
6:V:884:GLU:OE1	6:V:884:GLU:N	2.48	0.43
4:G:52:GLU:OE2	4:G:54:ARG:NE	2.51	0.42
6:Q:13:ARG:NH2	6:R:122:SER:O	2.51	0.42
4:E:210:ILE:HG21	4:E:218:ALA:HB2	2.00	0.42
6:T:568:GLU:OE1	6:T:862:ARG:NH1	2.52	0.42
4:E:185:VAL:HG23	4:E:190:VAL:CG2	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:R:568:GLU:OE1	6:R:862:ARG:NH1	2.52	0.42
1:A:48:GLU:N	1:A:48:GLU:OE1	2.52	0.42
4:F:544:ASP:OD1	4:F:545:PHE:N	2.52	0.42
4:G:38:LEU:HD21	4:G:415:VAL:HG13	2.01	0.42
4:G:226:LYS:HZ2	4:G:254:ASP:CG	2.27	0.42
6:Q:531:GLU:OE1	6:Q:531:GLU:N	2.46	0.42
2:C:253:ASP:OD1	2:C:254:VAL:N	2.52	0.42
6:V:1104:ASN:OD1	6:V:1107:PHE:N	2.51	0.42
4:F:34:LYS:NZ	4:F:495:GLU:OE2	2.36	0.42
5:N:6:TYR:CE2	5:N:99:LYS:HE2	2.54	0.42
6:U:810:GLY:N	6:U:937:GLY:O	2.52	0.42
6:V:531:GLU:OE1	6:V:531:GLU:N	2.46	0.42
4:I:167:PHE:HB3	4:I:241:ILE:HD11	2.02	0.42
4:J:505:LYS:NZ	4:J:509:GLU:OE2	2.52	0.42
6:Q:884:GLU:OE1	6:Q:884:GLU:N	2.48	0.42
4:F:153:ILE:HG23	4:F:154:PHE:CD1	2.54	0.42
4:H:244:LYS:O	4:H:246:VAL:HG23	2.20	0.42
5:O:134:LYS:HB3	6:V:8:SER:OG	2.18	0.42
6:V:295:PHE:CD2	6:V:295:PHE:N	2.87	0.42
4:G:172:ASP:O	4:G:176:GLN:N	2.52	0.41
6:T:94:ASP:OD1	6:T:95:SER:N	2.53	0.41
4:F:549:ASP:OD2	4:F:564:THR:OG1	2.33	0.41
6:R:94:ASP:OD1	6:R:95:SER:N	2.53	0.41
2:C:281:GLN:NE2	2:C:345:VAL:O	2.51	0.41
4:F:411:TYR:O	4:F:415:VAL:HG23	2.21	0.41
6:R:892:VAL:HG12	6:R:934:LEU:HD23	2.03	0.41
6:T:892:VAL:HG12	6:T:934:LEU:HD23	2.02	0.41
2:B:268:VAL:HG23	2:B:321:VAL:HG22	2.01	0.41
6:S:65:GLY:O	6:S:77:ASN:ND2	2.54	0.41
2:C:257:VAL:O	2:C:259:LYS:N	2.54	0.41
4:F:188:GLN:N	4:F:188:GLN:OE1	2.54	0.41
6:U:65:GLY:O	6:U:77:ASN:ND2	2.54	0.41
4:J:173:GLU:O	4:J:176:GLN:NE2	2.54	0.40
4:J:437:VAL:HG11	4:J:440:LEU:HD21	2.04	0.40
4:G:167:PHE:HB2	4:G:183:LEU:HD23	2.03	0.40
6:R:810:GLY:N	6:R:937:GLY:O	2.50	0.40
6:S:998:TRP:CD1	6:S:999:GLU:H	2.39	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	232/234 (99%)	223 (96%)	9 (4%)	0	100	100
2	B	346/348 (99%)	341 (99%)	5 (1%)	0	100	100
2	C	345/348 (99%)	339 (98%)	6 (2%)	0	100	100
3	D	367/1019 (36%)	354 (96%)	13 (4%)	0	100	100
4	E	534/587 (91%)	502 (94%)	31 (6%)	1 (0%)	43	77
4	F	547/587 (93%)	516 (94%)	31 (6%)	0	100	100
4	G	547/587 (93%)	521 (95%)	26 (5%)	0	100	100
4	H	547/587 (93%)	524 (96%)	23 (4%)	0	100	100
4	I	547/587 (93%)	526 (96%)	21 (4%)	0	100	100
4	J	547/587 (93%)	522 (95%)	25 (5%)	0	100	100
5	K	174/194 (90%)	166 (95%)	8 (5%)	0	100	100
5	L	174/194 (90%)	170 (98%)	4 (2%)	0	100	100
5	M	174/194 (90%)	165 (95%)	9 (5%)	0	100	100
5	N	174/194 (90%)	169 (97%)	5 (3%)	0	100	100
5	O	174/194 (90%)	169 (97%)	5 (3%)	0	100	100
5	P	174/194 (90%)	170 (98%)	4 (2%)	0	100	100
6	Q	900/1152 (78%)	876 (97%)	24 (3%)	0	100	100
6	R	900/1152 (78%)	872 (97%)	27 (3%)	1 (0%)	48	83
6	S	900/1152 (78%)	872 (97%)	28 (3%)	0	100	100
6	T	900/1152 (78%)	869 (97%)	30 (3%)	1 (0%)	48	83
6	U	898/1152 (78%)	866 (96%)	32 (4%)	0	100	100
6	V	900/1152 (78%)	862 (96%)	35 (4%)	3 (0%)	36	71
7	W	456/458 (100%)	441 (97%)	15 (3%)	0	100	100
7	X	456/458 (100%)	446 (98%)	10 (2%)	0	100	100
7	Y	456/458 (100%)	445 (98%)	11 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	12369/14921 (83%)	11926 (96%)	437 (4%)	6 (0%)	100	100

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	V	325	ASP
6	T	57	ASP
6	R	519	ILE
6	V	346	GLY
4	E	409	PRO
6	V	58	GLY

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	209/209 (100%)	208 (100%)	1 (0%)	81	81
2	B	311/311 (100%)	310 (100%)	1 (0%)	86	84
2	C	310/311 (100%)	309 (100%)	1 (0%)	86	84
3	D	162/928 (18%)	162 (100%)	0	100	100
4	E	459/495 (93%)	456 (99%)	3 (1%)	76	79
4	F	471/495 (95%)	464 (98%)	7 (2%)	57	70
4	G	471/495 (95%)	471 (100%)	0	100	100
4	H	471/495 (95%)	470 (100%)	1 (0%)	87	85
4	I	471/495 (95%)	464 (98%)	7 (2%)	57	70
4	J	471/495 (95%)	470 (100%)	1 (0%)	87	85
5	K	156/171 (91%)	154 (99%)	2 (1%)	61	72
5	L	156/171 (91%)	155 (99%)	1 (1%)	78	80
5	M	156/171 (91%)	155 (99%)	1 (1%)	78	80
5	N	156/171 (91%)	155 (99%)	1 (1%)	78	80
5	O	156/171 (91%)	155 (99%)	1 (1%)	78	80

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
5	P	156/171 (91%)	155 (99%)	1 (1%)	78	80
6	Q	800/1010 (79%)	797 (100%)	3 (0%)	84	82
6	R	800/1010 (79%)	800 (100%)	0	100	100
6	S	800/1010 (79%)	798 (100%)	2 (0%)	86	84
6	T	800/1010 (79%)	799 (100%)	1 (0%)	88	88
6	U	800/1010 (79%)	798 (100%)	2 (0%)	86	84
6	V	800/1010 (79%)	795 (99%)	5 (1%)	78	80
All	All	9542/11815 (81%)	9500 (100%)	42 (0%)	81	82

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ASP
2	B	271	SER
2	C	206	ASP
4	E	102	ILE
4	E	354	GLU
4	E	428	GLU
4	F	3	VAL
4	F	57	SER
4	F	114	ASN
4	F	247	TYR
4	F	453	ASN
4	F	481	ASN
4	F	482	ASP
4	H	453	ASN
4	I	48	ASN
4	I	141	ASP
4	I	167	PHE
4	I	241	ILE
4	I	247	TYR
4	I	264	ILE
4	I	557	ASN
4	J	202	TYR
5	K	172	GLU
5	K	176	ASN
5	L	176	ASN
5	M	168	ARG
5	N	176	ASN
5	O	176	ASN

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Mol	Chain	Res	Type
5	P	176	ASN
6	Q	33	GLN
6	Q	130	VAL
6	Q	1134	GLU
6	S	33	GLN
6	S	1133	THR
6	T	57	ASP
6	U	33	GLN
6	U	1133	THR
6	V	33	GLN
6	V	60	LYS
6	V	358	ARG
6	V	481	LEU
6	V	1134	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (69) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
2	B	111	GLN
4	E	269	GLN
4	E	342	GLN
4	E	376	GLN
4	E	543	GLN
4	F	407	HIS
4	F	497	ASN
4	F	543	GLN
4	G	126	ASN
4	G	269	GLN
4	I	120	GLN
4	I	171	HIS
4	I	204	ASN
4	I	258	GLN
4	I	262	ASN
4	I	540	ASN
4	J	120	GLN
4	J	204	ASN
4	J	262	ASN
4	J	465	ASN
5	K	45	ASN
5	K	50	GLN
5	L	82	ASN

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Mol	Chain	Res	Type
5	M	50	GLN
5	M	176	ASN
5	N	82	ASN
5	N	162	GLN
5	O	50	GLN
5	O	82	ASN
5	O	92	ASN
5	O	158	ASN
5	P	41	GLN
5	P	50	GLN
5	P	158	ASN
6	Q	293	GLN
6	Q	356	ASN
6	Q	822	ASN
6	R	293	GLN
6	R	356	ASN
6	R	822	ASN
6	R	910	ASN
6	S	339	GLN
6	S	341	GLN
6	T	4	ASN
6	T	61	GLN
6	T	168	GLN
6	T	248	ASN
6	T	293	GLN
6	T	339	GLN
6	T	441	ASN
6	T	488	ASN
6	T	822	ASN
6	T	910	ASN
6	T	1122	ASN
6	U	49	ASN
6	U	92	ASN
6	U	210	HIS
6	U	276	GLN
6	U	293	GLN
6	U	356	ASN
6	U	1105	ASN
6	V	77	ASN
6	V	195	ASN
6	V	248	ASN
6	V	276	GLN

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Mol	Chain	Res	Type
6	V	381	ASN
6	V	441	ASN
6	V	822	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.

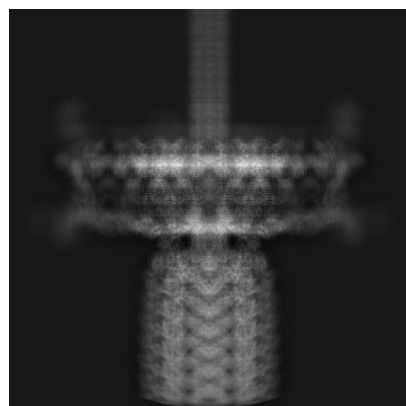
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-19971. These allow visual inspection of the internal detail of the map and identification of artifacts.

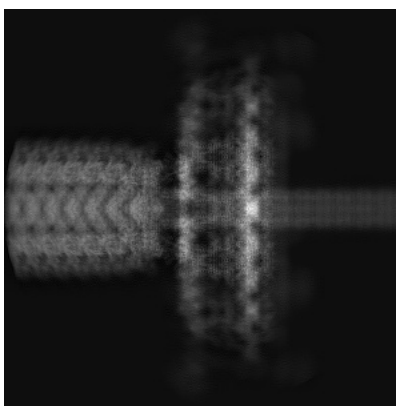
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

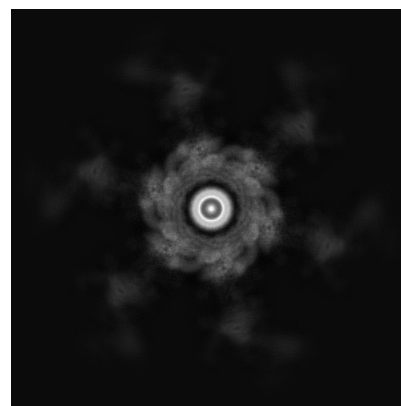
6.1.1 Primary map



X

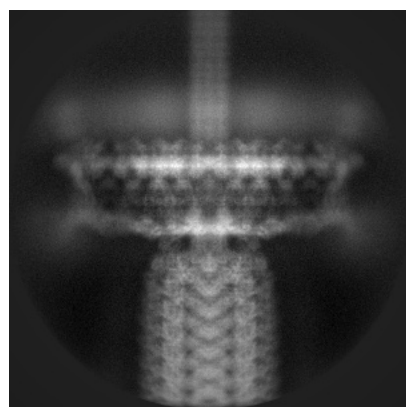


Y

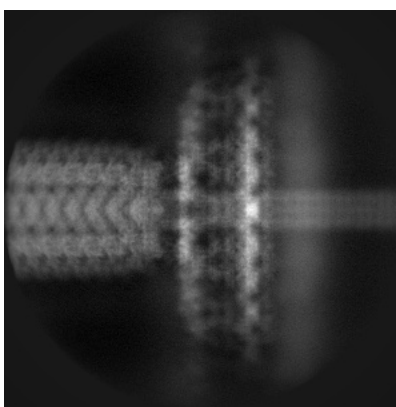


Z

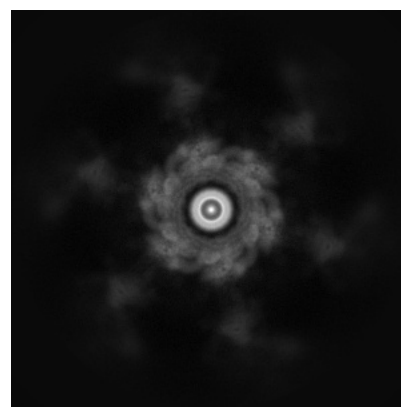
6.1.2 Raw map



X



Y

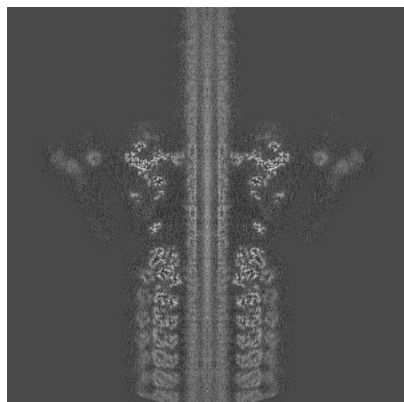


Z

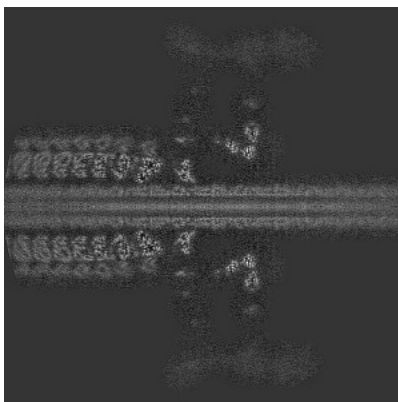
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

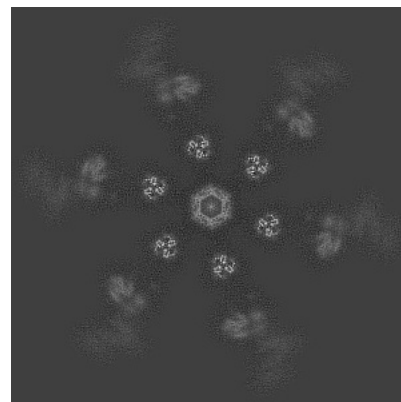
6.2.1 Primary map



X Index: 360

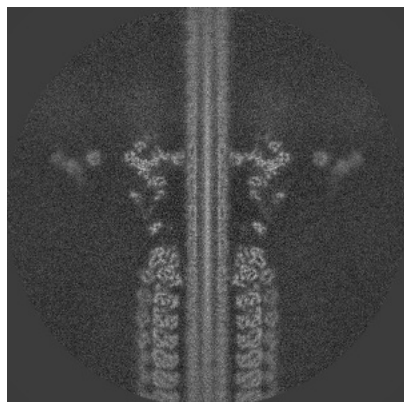


Y Index: 360

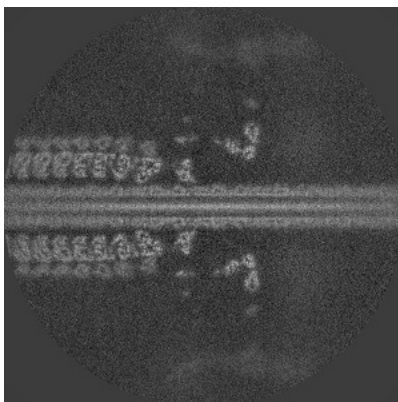


Z Index: 360

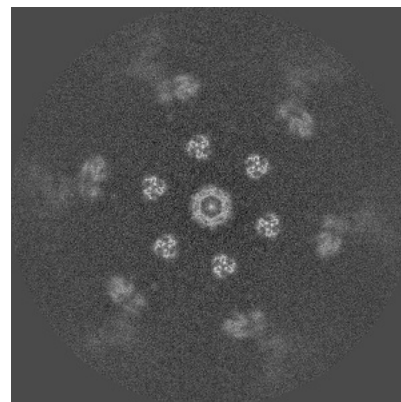
6.2.2 Raw map



X Index: 360



Y Index: 360



Z Index: 360

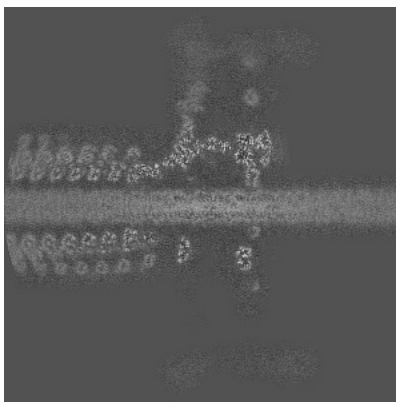
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

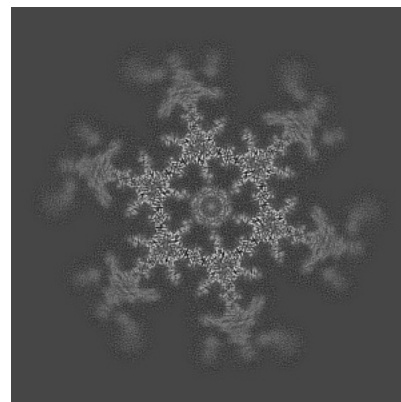
6.3.1 Primary map



X Index: 339

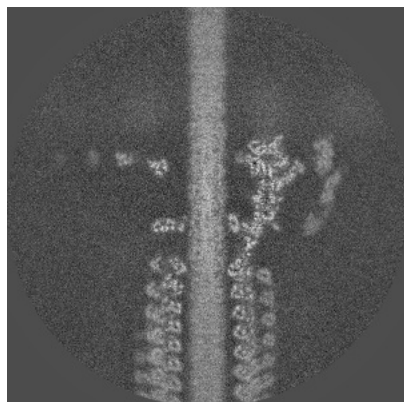


Y Index: 339

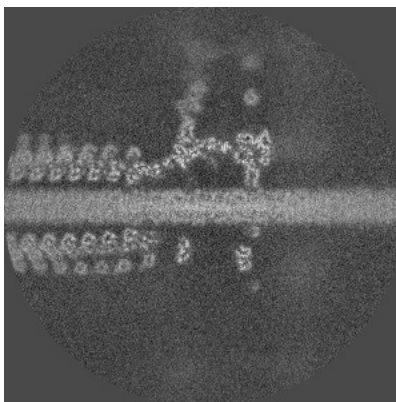


Z Index: 438

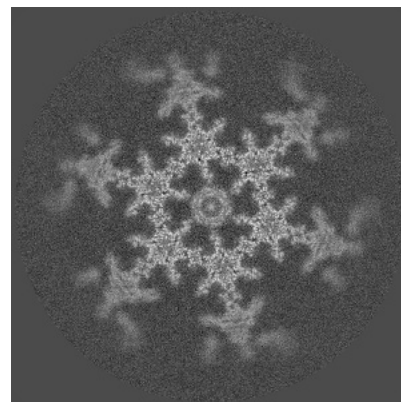
6.3.2 Raw map



X Index: 339



Y Index: 338

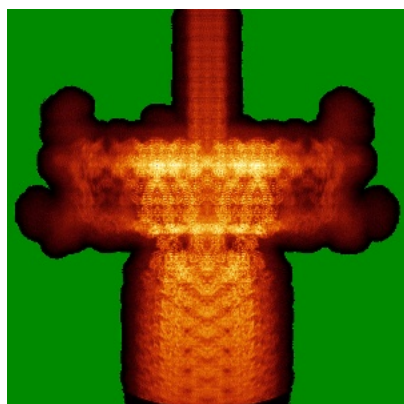


Z Index: 438

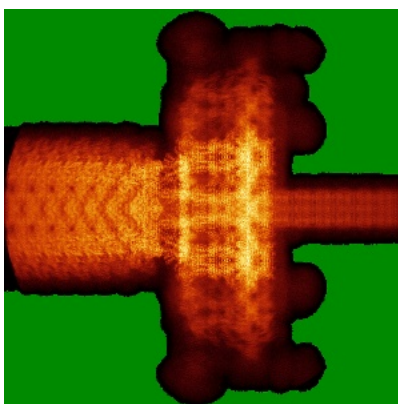
The images above show the largest variance slices of the map in three orthogonal directions.

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

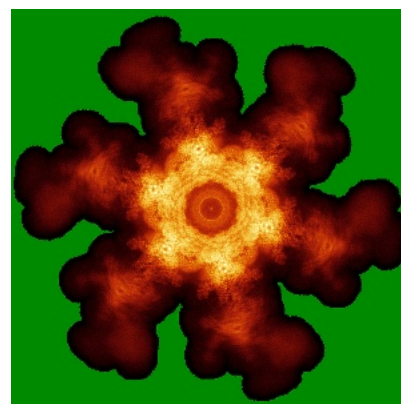
6.4.1 Primary map



X

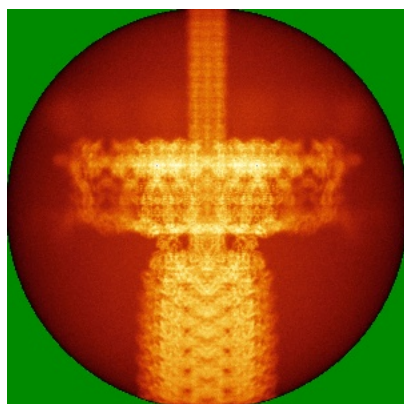


Y

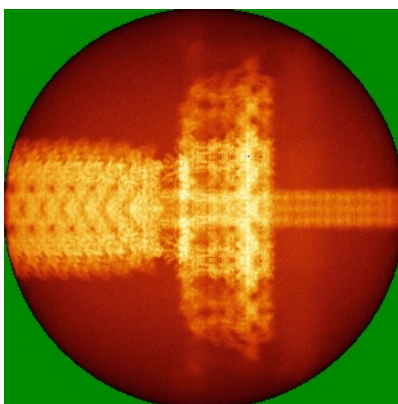


Z

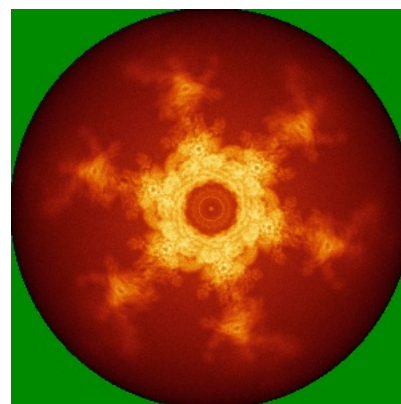
6.4.2 Raw map



X



Y

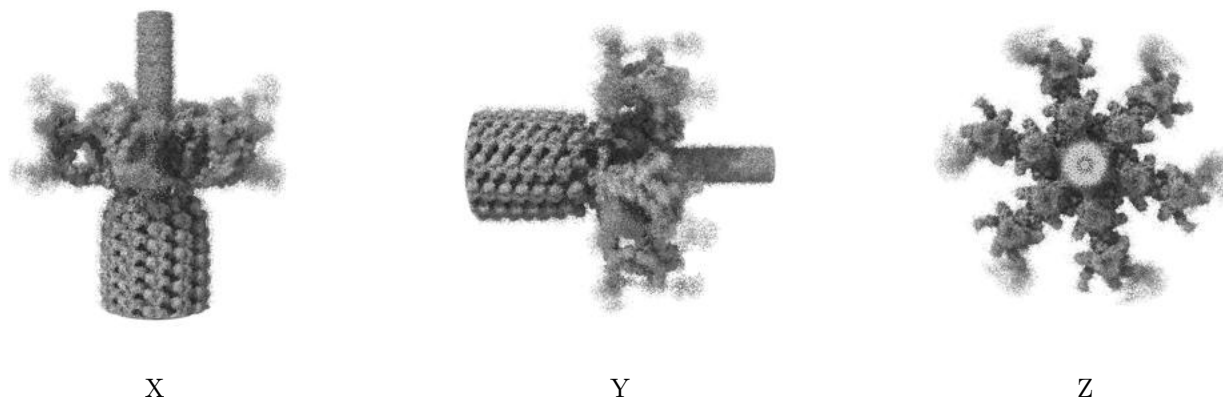


Z

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.

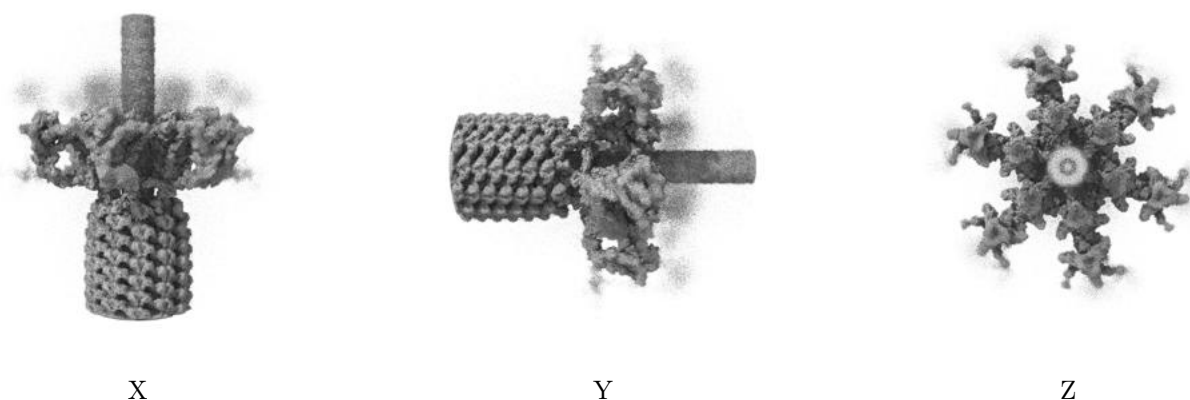
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.006. 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.

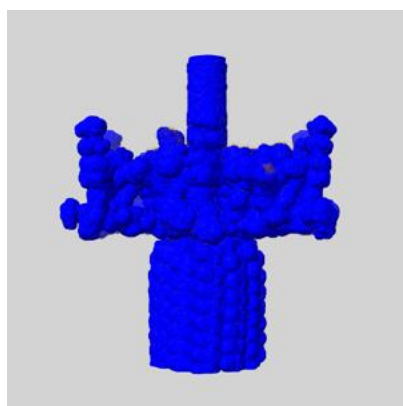
6.6 Mask visualisation [i](#)

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

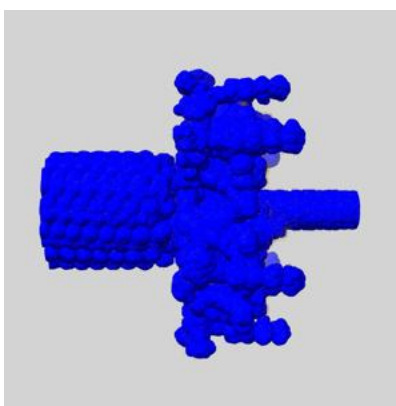
A mask typically either:

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

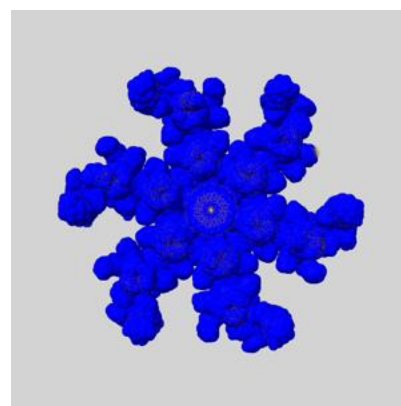
6.6.1 emd_19971_msk_1.map [i](#)



X



Y

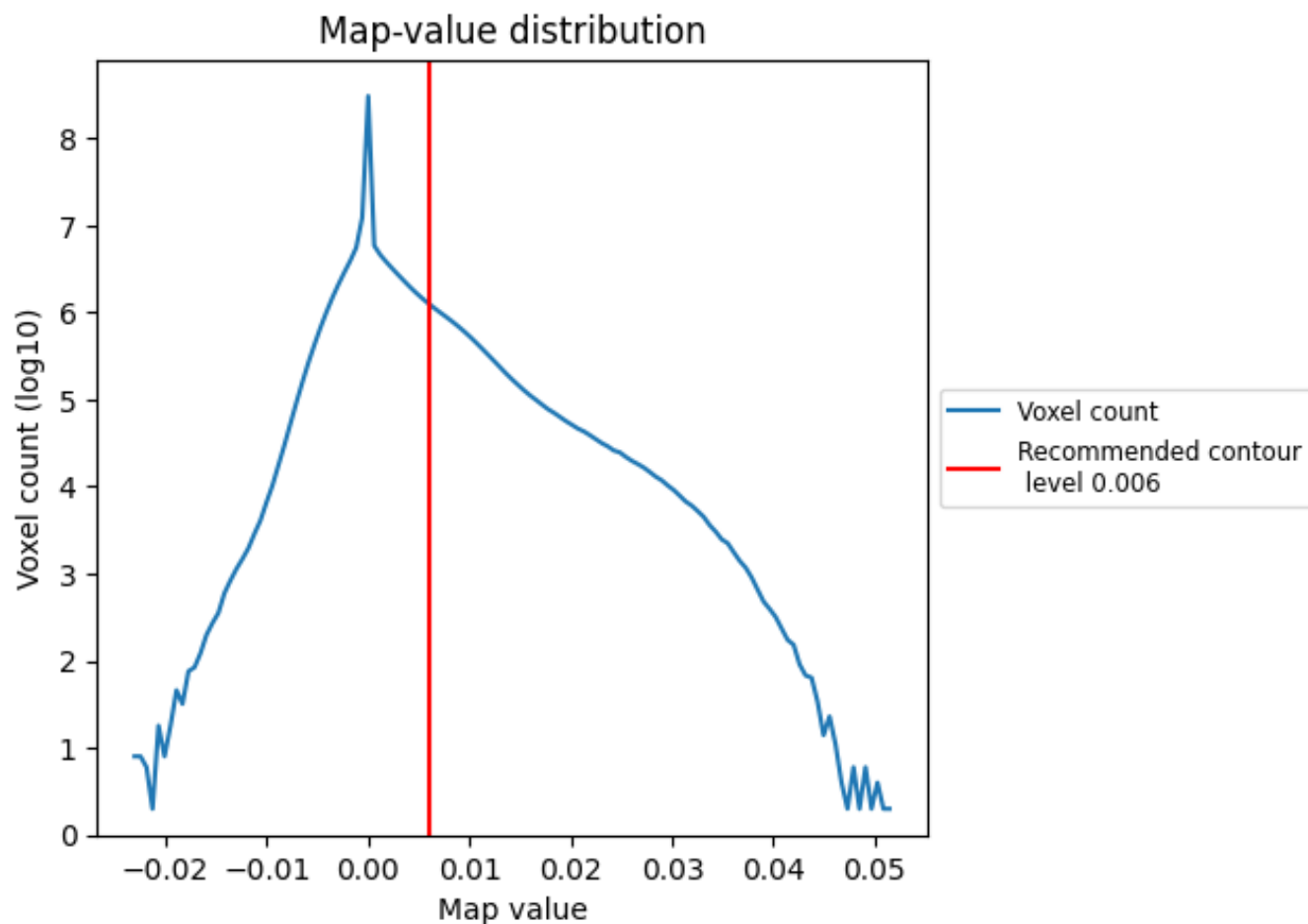


Z

7 Map analysis [i](#)

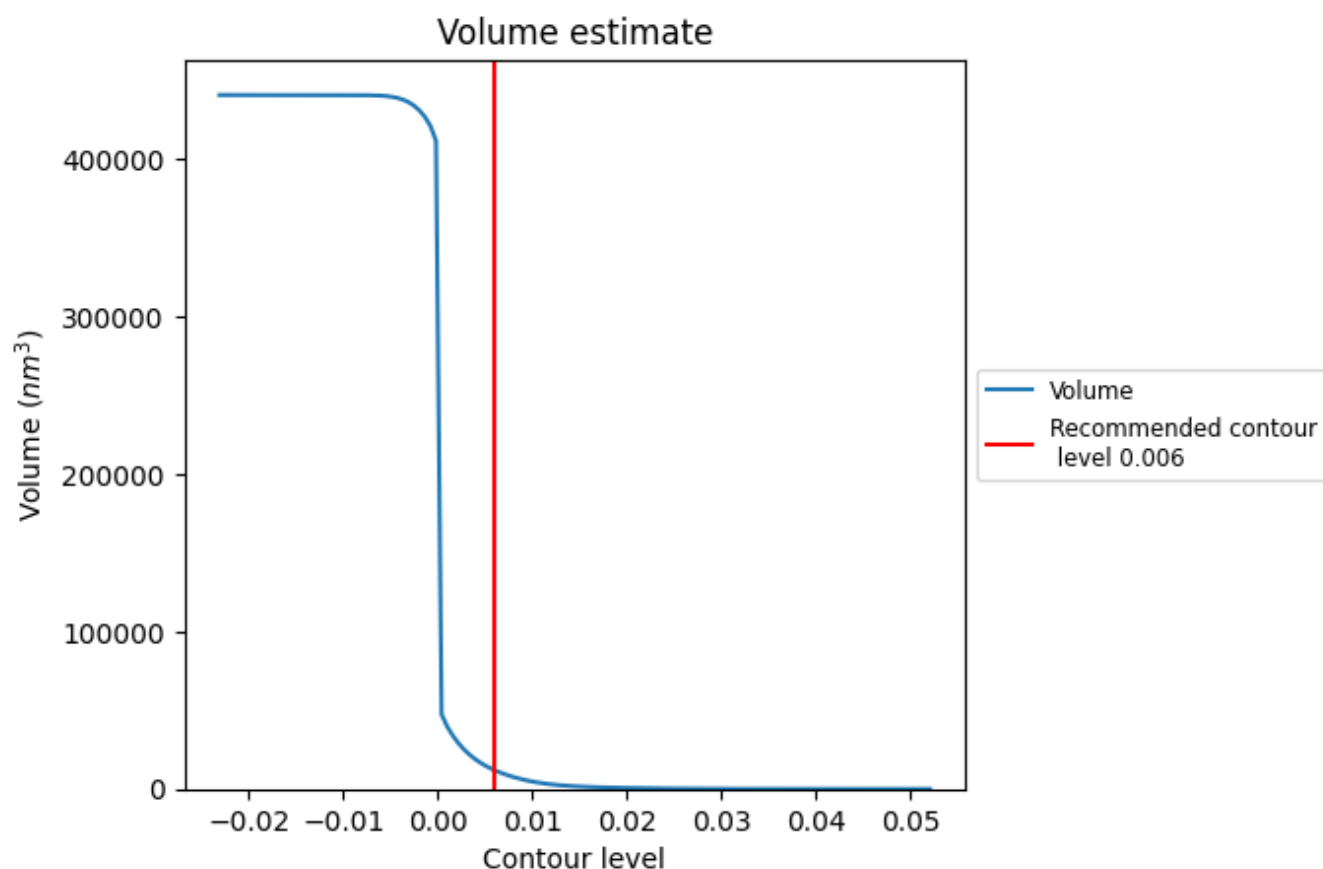
This section contains the results of statistical analysis of the map.

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.

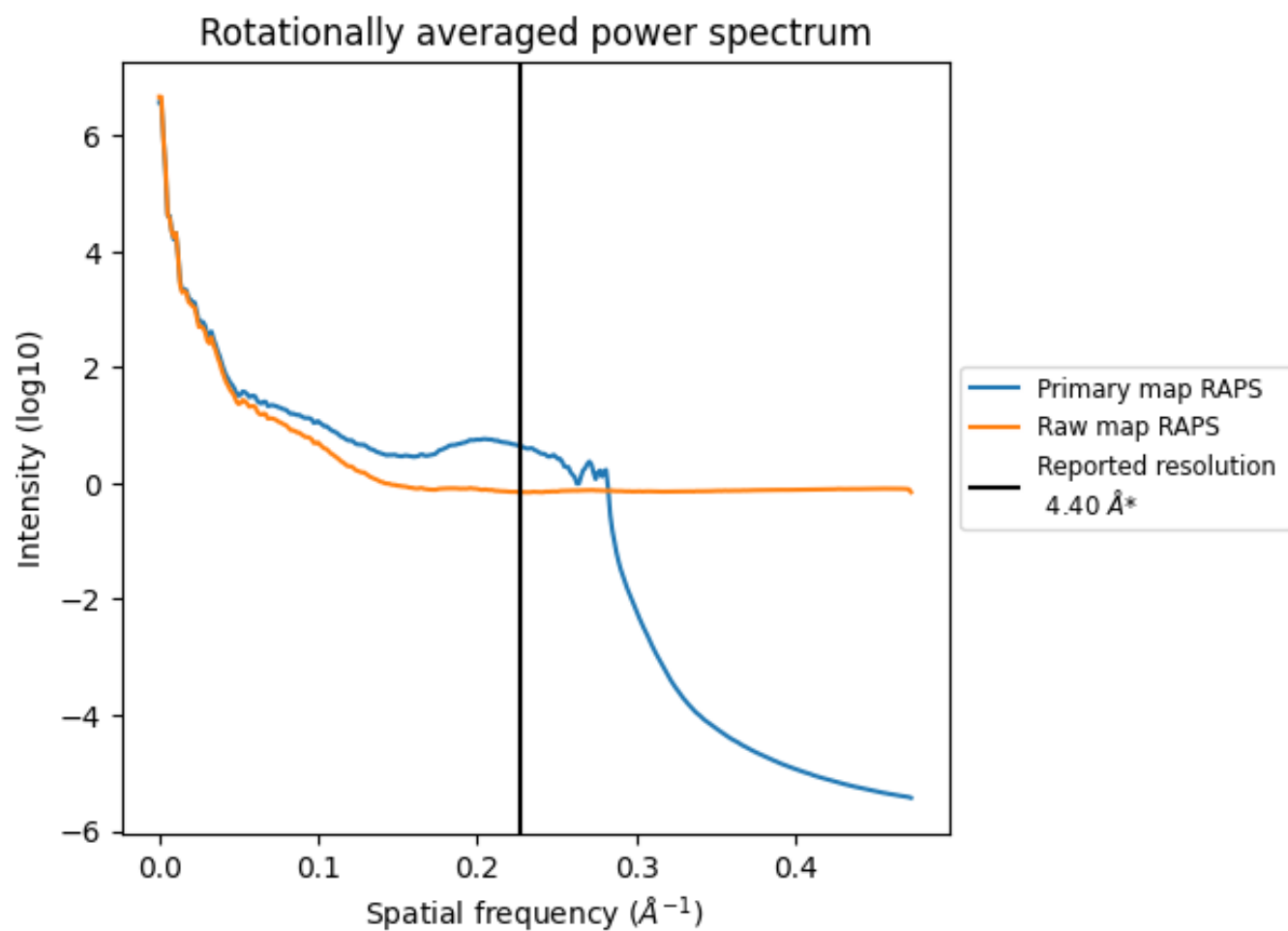
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 11953 nm^3 ; this corresponds to an approximate mass of 10798 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 ⓘ

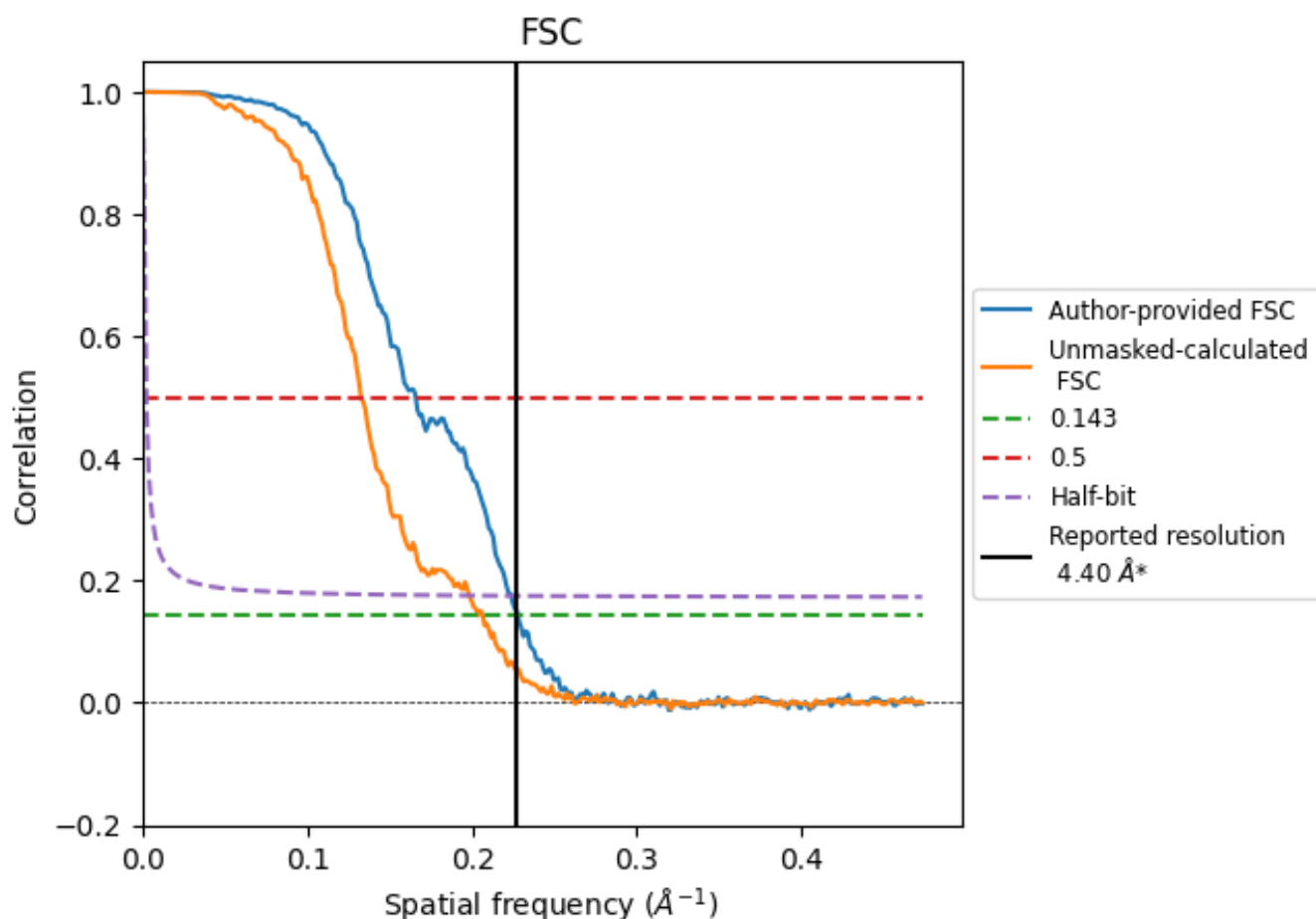


*Reported resolution corresponds to spatial frequency of 0.227 $\text{\AA}^{-1}$

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 $0.227 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.40	-	-
Author-provided FSC curve	4.39	6.03	4.47
Unmasked-calculated*	4.91	7.54	5.04

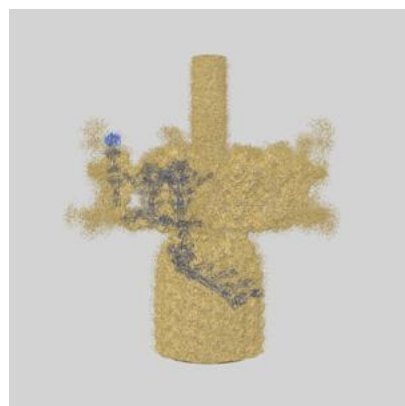
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.91 differs from the reported value 4.4 by more than 10 %

9 Map-model fit [i](#)

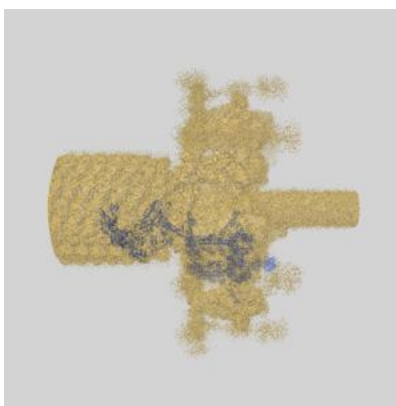
This section contains information regarding the fit between EMDB map EMD-19971 and PDB model 9EUI. Per-residue inclusion information can be found in section 3 on page 9.

9.1 Map-model overlays

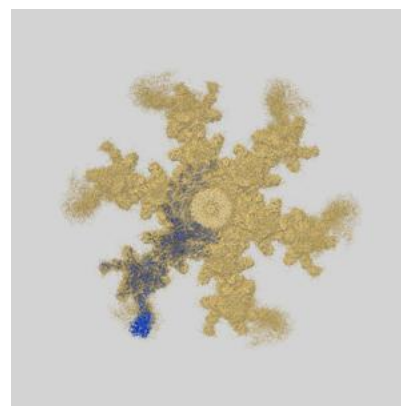
9.1.1 Map-model overlay [i](#)



X

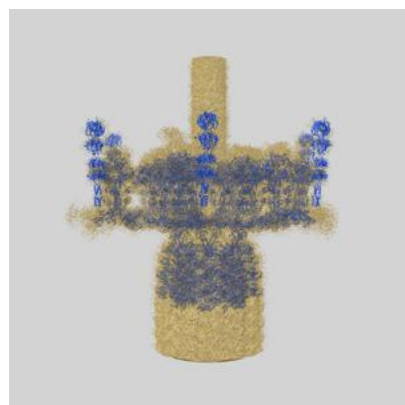


Y

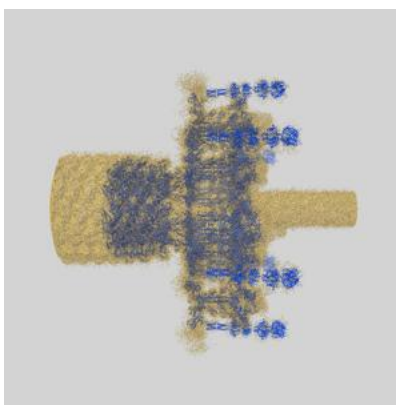


Z

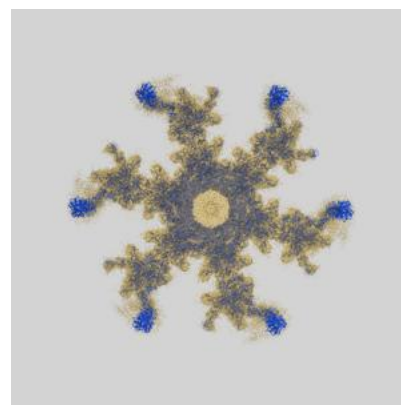
9.1.2 Map-model assembly overlay [i](#)



X



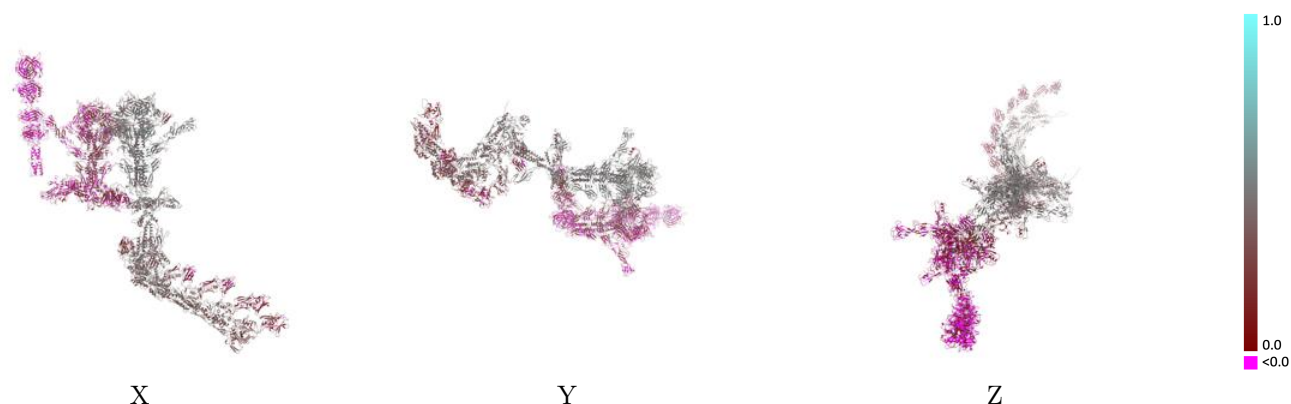
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.006 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](#)



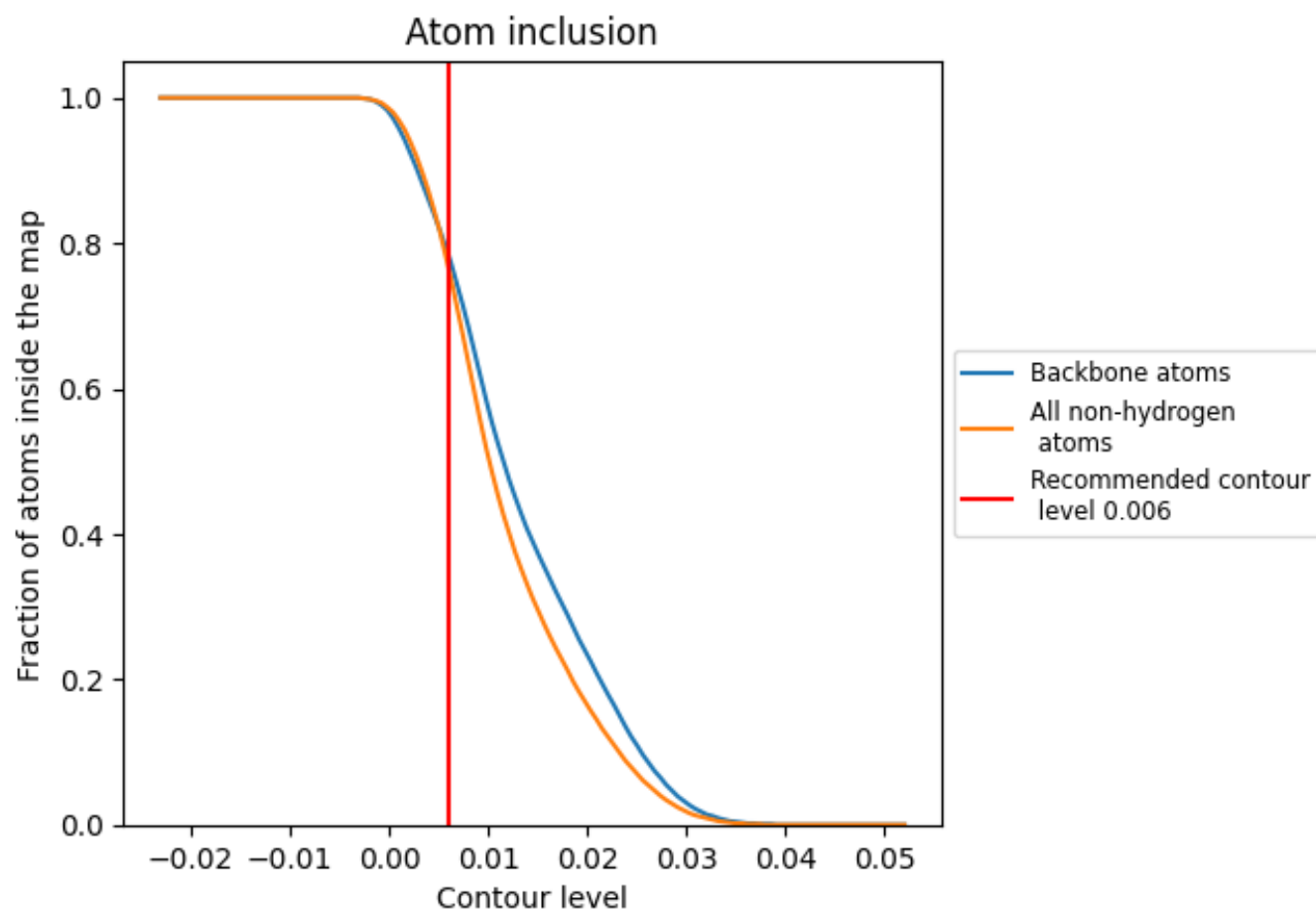
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.006).

9.4 Atom inclusion [i](#)



At the recommended contour level, 78% of all backbone atoms, 77% 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 (0.006) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7670	 0.2730
A	 0.9250	 0.4230
B	 0.9430	 0.4230
C	 0.9410	 0.4210
D	 0.8230	 0.3060
E	 0.8690	 0.3620
F	 0.9000	 0.3670
G	 0.8370	 0.3160
H	 0.8290	 0.3100
I	 0.8490	 0.3010
J	 0.8390	 0.2860
K	 0.8280	 0.1420
L	 0.8760	 0.2240
M	 0.6960	 0.1560
N	 0.4920	 0.1130
O	 0.1590	 0.0760
P	 0.0750	 0.0540
Q	 0.9440	 0.4360
R	 0.9430	 0.4360
S	 0.9440	 0.4300
T	 0.7550	 0.1600
U	 0.6990	 0.1010
V	 0.6770	 0.1130
W	 0.0250	 -0.0010
X	 0.0110	 -0.0060
Y	 0.0110	 -0.0030

