



wwPDB EM Validation Summary Report ⓘ

Mar 10, 2026 – 02:54 AM UTC

PDB ID : 5LZP / pdb_00005lzp
EMDB ID : EMD-4128
Title : Binding of the C-terminal GQYL motif of the bacterial proteasome activator Bpa to the 20S proteasome
Authors : Bolten, M.; Delley, C.L.; Leibundgut, M.; Boehringer, D.; Ban, N.; Weber-Ban, E.
Deposited on : 2016-09-30
Resolution : 3.50 Å(reported)

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:

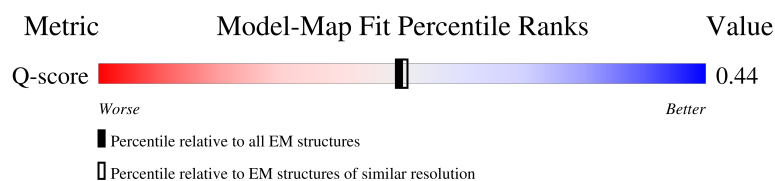
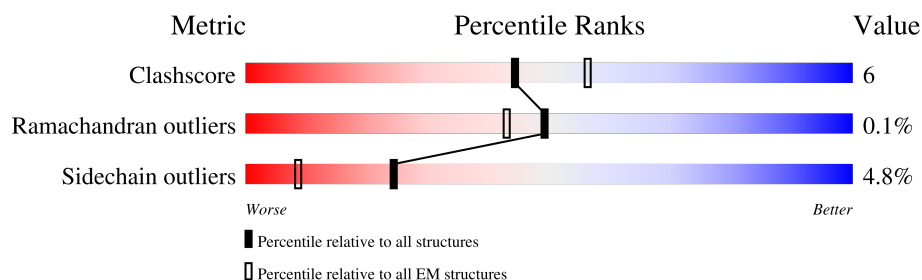
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.50 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.



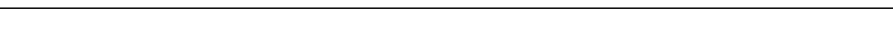
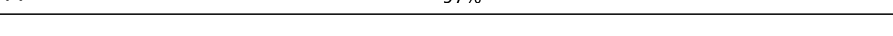
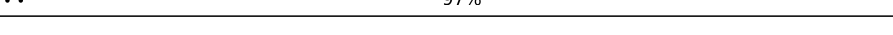
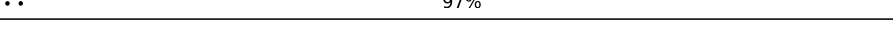
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13950 (3.00 - 4.00)

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	0	241	 70% 17% • 11%
1	2	241	 70% 17% • 11%
1	4	241	 70% 17% • 11%
1	6	241	 72% 15% • 11%

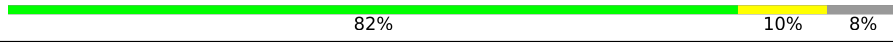
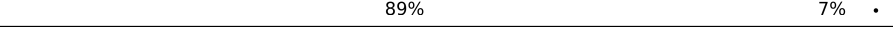
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Mol	Chain	Length	Quality of chain
1	8	241	
1	B	241	
1	D	241	
1	H	241	
1	J	241	
1	M	241	
1	Q	241	
1	S	241	
1	W	241	
1	Y	241	
2	1	180	
2	3	180	
2	5	180	
2	7	180	
2	V	180	
2	X	180	
2	Z	180	
3	A	242	
3	C	242	
3	E	242	
3	F	242	
3	G	242	
3	I	242	
3	K	242	
3	L	242	

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Mol	Chain	Length	Quality of chain
3	N	242	 83% 13% 8%
3	O	242	 82% 10% 8%
3	P	242	 89% 7% 8%
3	R	242	 79% 12% 8%
3	T	242	 79% 12% 8%
3	U	242	 85% 7% 8%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 46557 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 Proteasome subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	2	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	4	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	6	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	8	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	B	214	Total	C	N	O	S	0	0
			1650	1033	302	312	3		
1	D	217	Total	C	N	O	S	0	0
			1668	1044	305	316	3		
1	H	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	J	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	M	214	Total	C	N	O	S	0	0
			1650	1033	302	312	3		
1	Q	214	Total	C	N	O	S	0	0
			1650	1033	302	312	3		
1	S	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	W	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
1	Y	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		

- Molecule 2 is a protein called Bacterial proteasome activator.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	1	5	Total	C	N	O	0	0
			39	25	6	8		

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Mol	Chain	Residues	Atoms				AltConf	Trace
2	3	5	Total	C	N	O	0	0
			39	25	6	8		
2	5	5	Total	C	N	O	0	0
			39	25	6	8		
2	7	5	Total	C	N	O	0	0
			39	25	6	8		
2	V	5	Total	C	N	O	0	0
			39	25	6	8		
2	X	5	Total	C	N	O	0	0
			39	25	6	8		
2	Z	5	Total	C	N	O	0	0
			39	25	6	8		

There are 49 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	-5	MET	-	initiating methionine	UNP P9WKX3
1	-4	HIS	-	expression tag	UNP P9WKX3
1	-3	HIS	-	expression tag	UNP P9WKX3
1	-2	HIS	-	expression tag	UNP P9WKX3
1	-1	HIS	-	expression tag	UNP P9WKX3
1	0	HIS	-	expression tag	UNP P9WKX3
1	1	HIS	-	expression tag	UNP P9WKX3
3	-5	MET	-	initiating methionine	UNP P9WKX3
3	-4	HIS	-	expression tag	UNP P9WKX3
3	-3	HIS	-	expression tag	UNP P9WKX3
3	-2	HIS	-	expression tag	UNP P9WKX3
3	-1	HIS	-	expression tag	UNP P9WKX3
3	0	HIS	-	expression tag	UNP P9WKX3
3	1	HIS	-	expression tag	UNP P9WKX3
5	-5	MET	-	initiating methionine	UNP P9WKX3
5	-4	HIS	-	expression tag	UNP P9WKX3
5	-3	HIS	-	expression tag	UNP P9WKX3
5	-2	HIS	-	expression tag	UNP P9WKX3
5	-1	HIS	-	expression tag	UNP P9WKX3
5	0	HIS	-	expression tag	UNP P9WKX3
5	1	HIS	-	expression tag	UNP P9WKX3
7	-5	MET	-	initiating methionine	UNP P9WKX3
7	-4	HIS	-	expression tag	UNP P9WKX3
7	-3	HIS	-	expression tag	UNP P9WKX3
7	-2	HIS	-	expression tag	UNP P9WKX3
7	-1	HIS	-	expression tag	UNP P9WKX3
7	0	HIS	-	expression tag	UNP P9WKX3

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Chain	Residue	Modelled	Actual	Comment	Reference
7	1	HIS	-	expression tag	UNP P9WKX3
V	-5	MET	-	initiating methionine	UNP P9WKX3
V	-4	HIS	-	expression tag	UNP P9WKX3
V	-3	HIS	-	expression tag	UNP P9WKX3
V	-2	HIS	-	expression tag	UNP P9WKX3
V	-1	HIS	-	expression tag	UNP P9WKX3
V	0	HIS	-	expression tag	UNP P9WKX3
V	1	HIS	-	expression tag	UNP P9WKX3
X	-5	MET	-	initiating methionine	UNP P9WKX3
X	-4	HIS	-	expression tag	UNP P9WKX3
X	-3	HIS	-	expression tag	UNP P9WKX3
X	-2	HIS	-	expression tag	UNP P9WKX3
X	-1	HIS	-	expression tag	UNP P9WKX3
X	0	HIS	-	expression tag	UNP P9WKX3
X	1	HIS	-	expression tag	UNP P9WKX3
Z	-5	MET	-	initiating methionine	UNP P9WKX3
Z	-4	HIS	-	expression tag	UNP P9WKX3
Z	-3	HIS	-	expression tag	UNP P9WKX3
Z	-2	HIS	-	expression tag	UNP P9WKX3
Z	-1	HIS	-	expression tag	UNP P9WKX3
Z	0	HIS	-	expression tag	UNP P9WKX3
Z	1	HIS	-	expression tag	UNP P9WKX3

- Molecule 3 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	A	222	Total	C	N	O	S	0	0
			1636	1026	282	323	5		
3	C	223	Total	C	N	O	S	0	0
			1640	1028	283	324	5		
3	E	222	Total	C	N	O	S	0	0
			1636	1026	282	323	5		
3	F	222	Total	C	N	O	S	0	0
			1636	1026	282	323	5		
3	G	222	Total	C	N	O	S	0	0
			1636	1026	282	323	5		
3	I	223	Total	C	N	O	S	0	0
			1640	1028	283	324	5		
3	K	222	Total	C	N	O	S	0	0
			1636	1026	282	323	5		
3	L	222	Total	C	N	O	S	0	0
			1636	1026	282	323	5		

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Mol	Chain	Residues	Atoms					AltConf	Trace
3	N	234	Total	C	N	O	S	0	0
			1715	1072	295	343	5		
3	O	223	Total	C	N	O	S	0	0
			1640	1028	283	324	5		
3	P	234	Total	C	N	O	S	0	0
			1715	1072	295	343	5		
3	R	223	Total	C	N	O	S	0	0
			1640	1028	283	324	5		
3	T	223	Total	C	N	O	S	0	0
			1640	1028	283	324	5		
3	U	223	Total	C	N	O	S	0	0
			1640	1028	283	324	5		

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	301	ALA	-	expression tag	UNP P9WHT9
A	535	TRP	-	expression tag	UNP P9WHT9
A	536	SER	-	expression tag	UNP P9WHT9
A	537	HIS	-	expression tag	UNP P9WHT9
A	538	PRO	-	expression tag	UNP P9WHT9
A	539	GLN	-	expression tag	UNP P9WHT9
A	540	PHE	-	expression tag	UNP P9WHT9
A	541	GLU	-	expression tag	UNP P9WHT9
A	542	LYS	-	expression tag	UNP P9WHT9
C	301	ALA	-	expression tag	UNP P9WHT9
C	535	TRP	-	expression tag	UNP P9WHT9
C	536	SER	-	expression tag	UNP P9WHT9
C	537	HIS	-	expression tag	UNP P9WHT9
C	538	PRO	-	expression tag	UNP P9WHT9
C	539	GLN	-	expression tag	UNP P9WHT9
C	540	PHE	-	expression tag	UNP P9WHT9
C	541	GLU	-	expression tag	UNP P9WHT9
C	542	LYS	-	expression tag	UNP P9WHT9
E	301	ALA	-	expression tag	UNP P9WHT9
E	535	TRP	-	expression tag	UNP P9WHT9
E	536	SER	-	expression tag	UNP P9WHT9
E	537	HIS	-	expression tag	UNP P9WHT9
E	538	PRO	-	expression tag	UNP P9WHT9
E	539	GLN	-	expression tag	UNP P9WHT9
E	540	PHE	-	expression tag	UNP P9WHT9
E	541	GLU	-	expression tag	UNP P9WHT9
E	542	LYS	-	expression tag	UNP P9WHT9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	301	ALA	-	expression tag	UNP P9WHT9
F	535	TRP	-	expression tag	UNP P9WHT9
F	536	SER	-	expression tag	UNP P9WHT9
F	537	HIS	-	expression tag	UNP P9WHT9
F	538	PRO	-	expression tag	UNP P9WHT9
F	539	GLN	-	expression tag	UNP P9WHT9
F	540	PHE	-	expression tag	UNP P9WHT9
F	541	GLU	-	expression tag	UNP P9WHT9
F	542	LYS	-	expression tag	UNP P9WHT9
G	301	ALA	-	expression tag	UNP P9WHT9
G	535	TRP	-	expression tag	UNP P9WHT9
G	536	SER	-	expression tag	UNP P9WHT9
G	537	HIS	-	expression tag	UNP P9WHT9
G	538	PRO	-	expression tag	UNP P9WHT9
G	539	GLN	-	expression tag	UNP P9WHT9
G	540	PHE	-	expression tag	UNP P9WHT9
G	541	GLU	-	expression tag	UNP P9WHT9
G	542	LYS	-	expression tag	UNP P9WHT9
I	301	ALA	-	expression tag	UNP P9WHT9
I	535	TRP	-	expression tag	UNP P9WHT9
I	536	SER	-	expression tag	UNP P9WHT9
I	537	HIS	-	expression tag	UNP P9WHT9
I	538	PRO	-	expression tag	UNP P9WHT9
I	539	GLN	-	expression tag	UNP P9WHT9
I	540	PHE	-	expression tag	UNP P9WHT9
I	541	GLU	-	expression tag	UNP P9WHT9
I	542	LYS	-	expression tag	UNP P9WHT9
K	301	ALA	-	expression tag	UNP P9WHT9
K	535	TRP	-	expression tag	UNP P9WHT9
K	536	SER	-	expression tag	UNP P9WHT9
K	537	HIS	-	expression tag	UNP P9WHT9
K	538	PRO	-	expression tag	UNP P9WHT9
K	539	GLN	-	expression tag	UNP P9WHT9
K	540	PHE	-	expression tag	UNP P9WHT9
K	541	GLU	-	expression tag	UNP P9WHT9
K	542	LYS	-	expression tag	UNP P9WHT9
L	301	ALA	-	expression tag	UNP P9WHT9
L	535	TRP	-	expression tag	UNP P9WHT9
L	536	SER	-	expression tag	UNP P9WHT9
L	537	HIS	-	expression tag	UNP P9WHT9
L	538	PRO	-	expression tag	UNP P9WHT9
L	539	GLN	-	expression tag	UNP P9WHT9

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Chain	Residue	Modelled	Actual	Comment	Reference
L	540	PHE	-	expression tag	UNP P9WHT9
L	541	GLU	-	expression tag	UNP P9WHT9
L	542	LYS	-	expression tag	UNP P9WHT9
N	301	ALA	-	expression tag	UNP P9WHT9
N	535	TRP	-	expression tag	UNP P9WHT9
N	536	SER	-	expression tag	UNP P9WHT9
N	537	HIS	-	expression tag	UNP P9WHT9
N	538	PRO	-	expression tag	UNP P9WHT9
N	539	GLN	-	expression tag	UNP P9WHT9
N	540	PHE	-	expression tag	UNP P9WHT9
N	541	GLU	-	expression tag	UNP P9WHT9
N	542	LYS	-	expression tag	UNP P9WHT9
O	301	ALA	-	expression tag	UNP P9WHT9
O	535	TRP	-	expression tag	UNP P9WHT9
O	536	SER	-	expression tag	UNP P9WHT9
O	537	HIS	-	expression tag	UNP P9WHT9
O	538	PRO	-	expression tag	UNP P9WHT9
O	539	GLN	-	expression tag	UNP P9WHT9
O	540	PHE	-	expression tag	UNP P9WHT9
O	541	GLU	-	expression tag	UNP P9WHT9
O	542	LYS	-	expression tag	UNP P9WHT9
P	301	ALA	-	expression tag	UNP P9WHT9
P	535	TRP	-	expression tag	UNP P9WHT9
P	536	SER	-	expression tag	UNP P9WHT9
P	537	HIS	-	expression tag	UNP P9WHT9
P	538	PRO	-	expression tag	UNP P9WHT9
P	539	GLN	-	expression tag	UNP P9WHT9
P	540	PHE	-	expression tag	UNP P9WHT9
P	541	GLU	-	expression tag	UNP P9WHT9
P	542	LYS	-	expression tag	UNP P9WHT9
R	301	ALA	-	expression tag	UNP P9WHT9
R	535	TRP	-	expression tag	UNP P9WHT9
R	536	SER	-	expression tag	UNP P9WHT9
R	537	HIS	-	expression tag	UNP P9WHT9
R	538	PRO	-	expression tag	UNP P9WHT9
R	539	GLN	-	expression tag	UNP P9WHT9
R	540	PHE	-	expression tag	UNP P9WHT9
R	541	GLU	-	expression tag	UNP P9WHT9
R	542	LYS	-	expression tag	UNP P9WHT9
T	301	ALA	-	expression tag	UNP P9WHT9
T	535	TRP	-	expression tag	UNP P9WHT9
T	536	SER	-	expression tag	UNP P9WHT9

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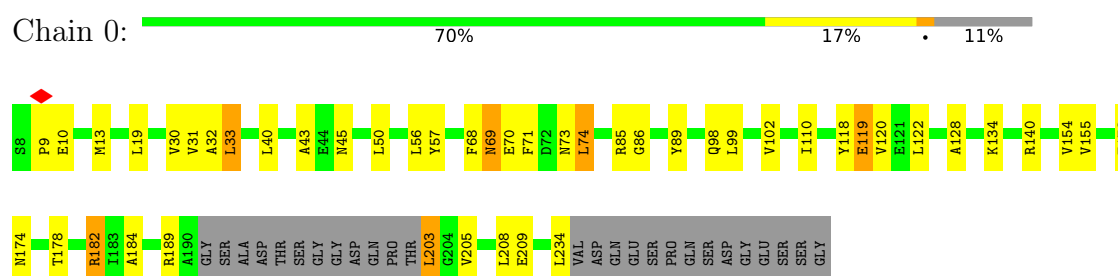
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Chain	Residue	Modelled	Actual	Comment	Reference
T	537	HIS	-	expression tag	UNP P9WHT9
T	538	PRO	-	expression tag	UNP P9WHT9
T	539	GLN	-	expression tag	UNP P9WHT9
T	540	PHE	-	expression tag	UNP P9WHT9
T	541	GLU	-	expression tag	UNP P9WHT9
T	542	LYS	-	expression tag	UNP P9WHT9
U	301	ALA	-	expression tag	UNP P9WHT9
U	535	TRP	-	expression tag	UNP P9WHT9
U	536	SER	-	expression tag	UNP P9WHT9
U	537	HIS	-	expression tag	UNP P9WHT9
U	538	PRO	-	expression tag	UNP P9WHT9
U	539	GLN	-	expression tag	UNP P9WHT9
U	540	PHE	-	expression tag	UNP P9WHT9
U	541	GLU	-	expression tag	UNP P9WHT9
U	542	LYS	-	expression tag	UNP P9WHT9

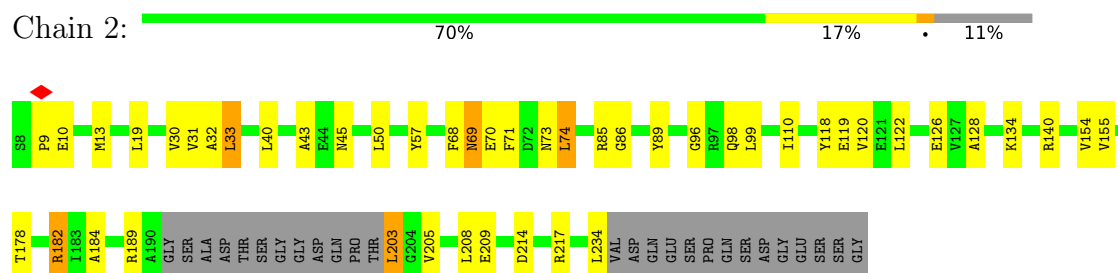
3 Residue-property plots

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.

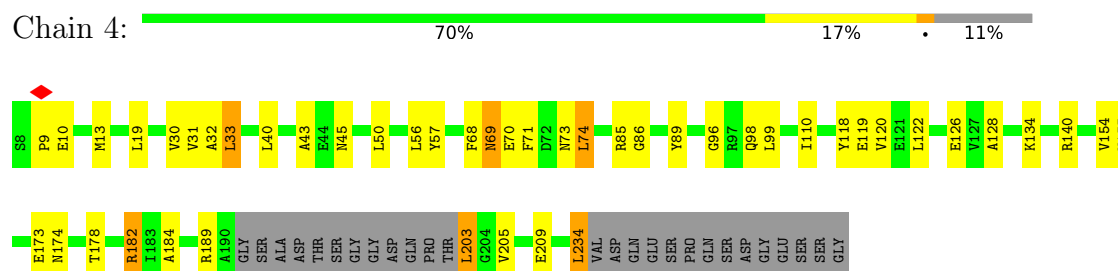
- Molecule 1: Proteasome subunit alpha



- Molecule 1: Proteasome subunit alpha

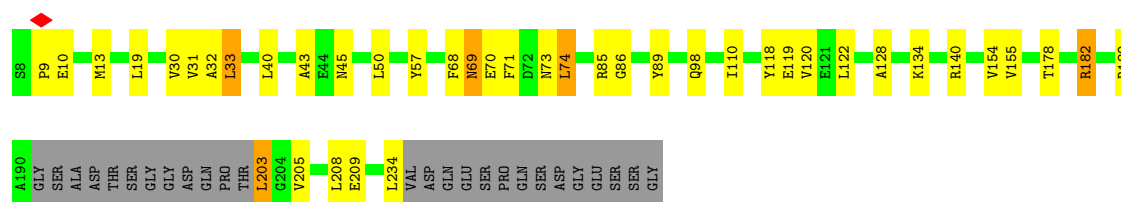


- Molecule 1: Proteasome subunit alpha

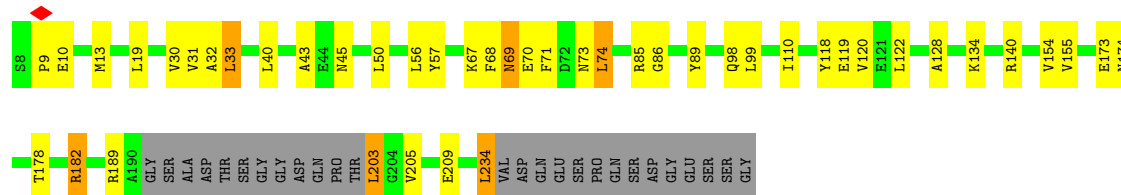


- Molecule 1: Proteasome subunit alpha

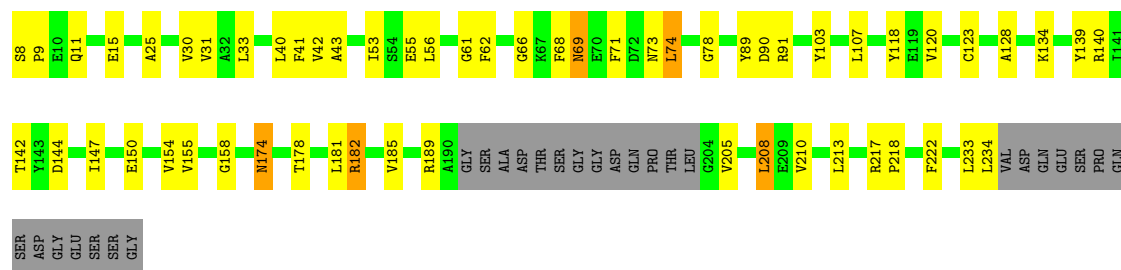




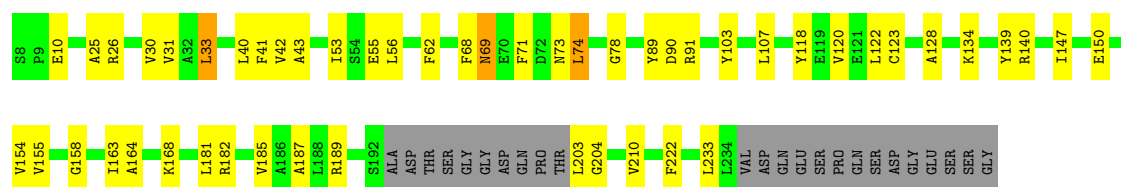
- Molecule 1: Proteasome subunit alpha



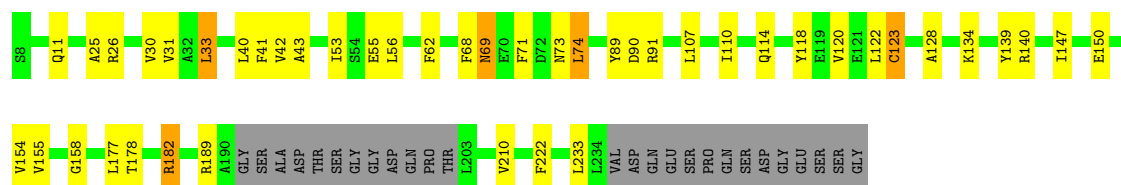
- Molecule 1: Proteasome subunit alpha



- Molecule 1: Proteasome subunit alpha

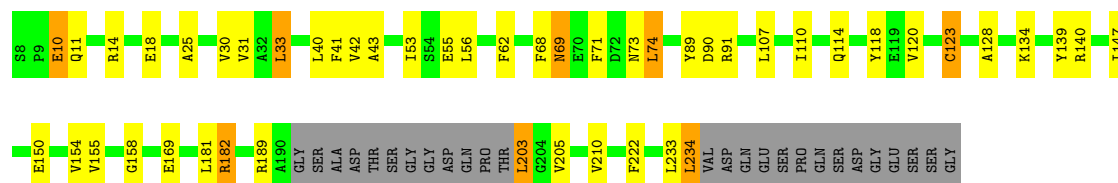


- Molecule 1: Proteasome subunit alpha



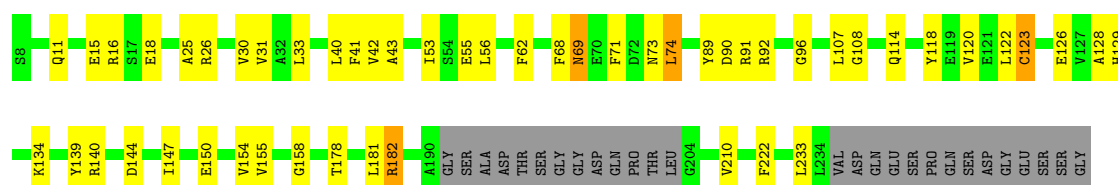
- Molecule 1: Proteasome subunit alpha

Chain J:  69% 17% 11%



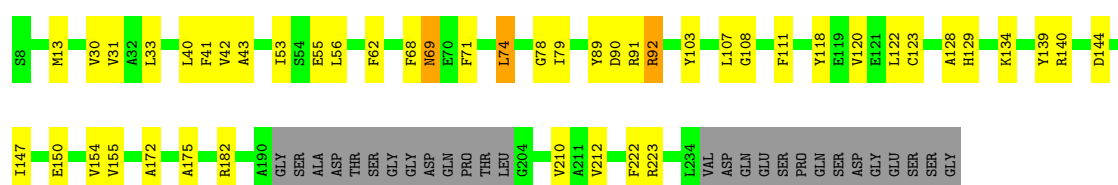
- Molecule 1: Proteasome subunit alpha

Chain M:  67% 20% 11%



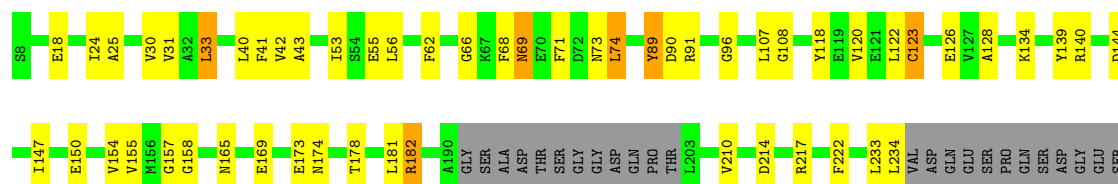
- Molecule 1: Proteasome subunit alpha

Chain Q:  69% 18% 11%



- Molecule 1: Proteasome subunit alpha

Chain S:  66% 20% 11%



SER
GLY

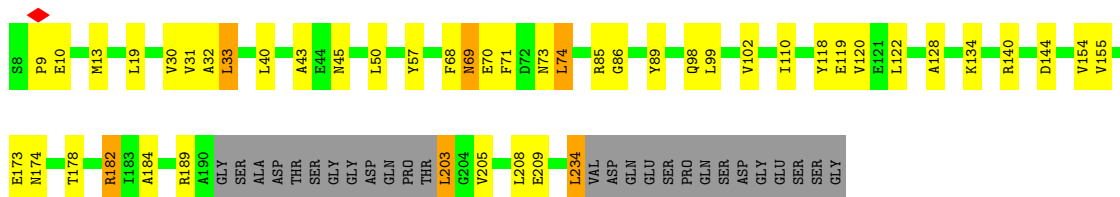
- Molecule 1: Proteasome subunit alpha

Chain W:  71% 16% 11%



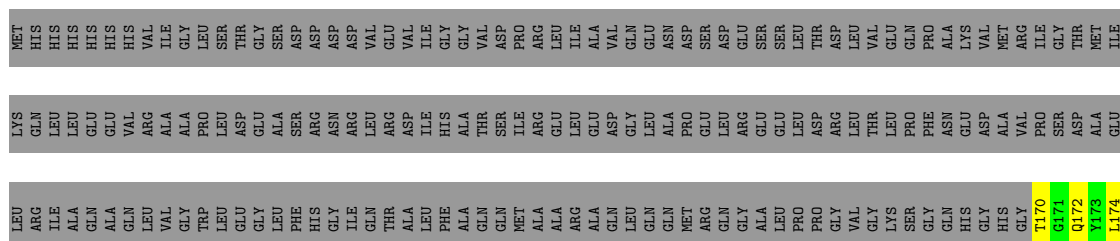
- Molecule 1: Proteasome subunit alpha

Chain Y: 70% 17% 1% 11%



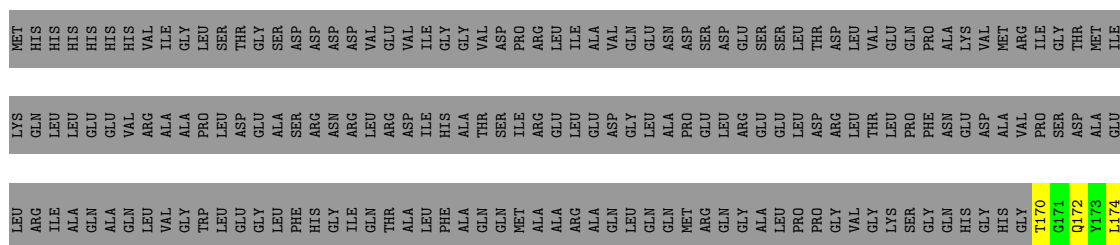
- Molecule 2: Bacterial proteasome activator

Chain 1: 97%

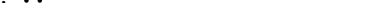


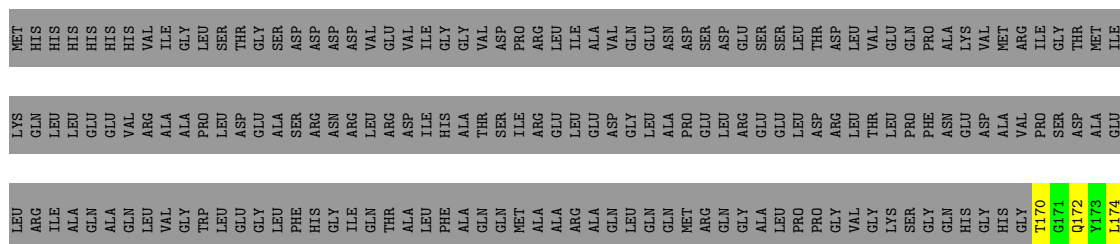
- Molecule 2: Bacterial proteasome activator

Chain 3: .. 97%



- Molecule 2: Bacterial proteasome activator

Chain 5:  97%



- Molecule 2: Bacterial proteasome activator

97%

- Molecule 2: Bacterial proteasome activator

97%

- Molecule 2: Bacterial proteasome activator

97%

- Molecule 2: Bacterial proteasome activator

97%

- Molecule 3: Proteasome subunit beta

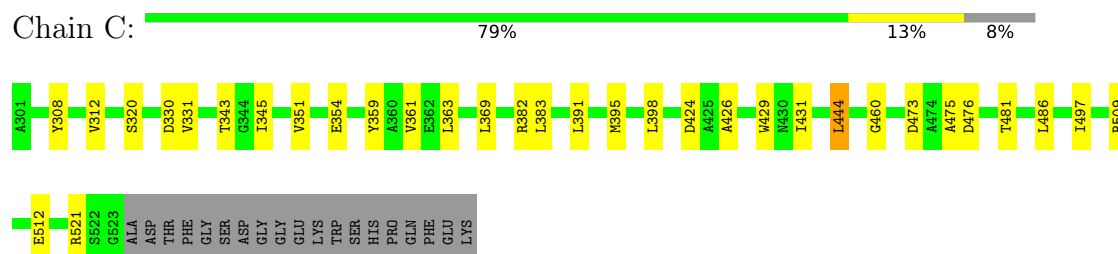
83%

7%

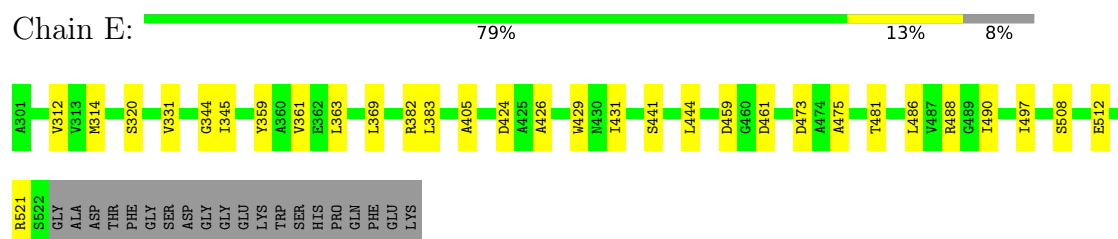
8%



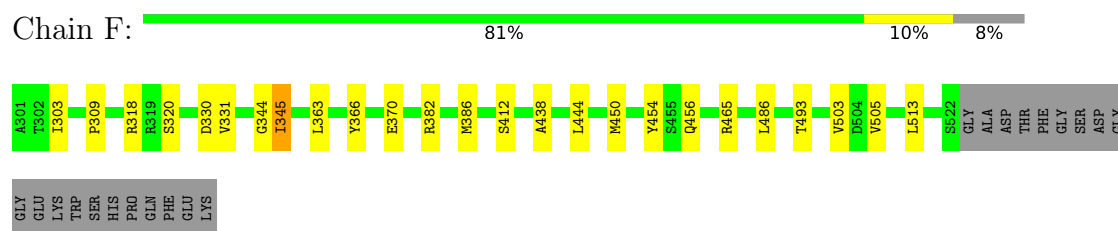
● Molecule 3: Proteasome subunit beta



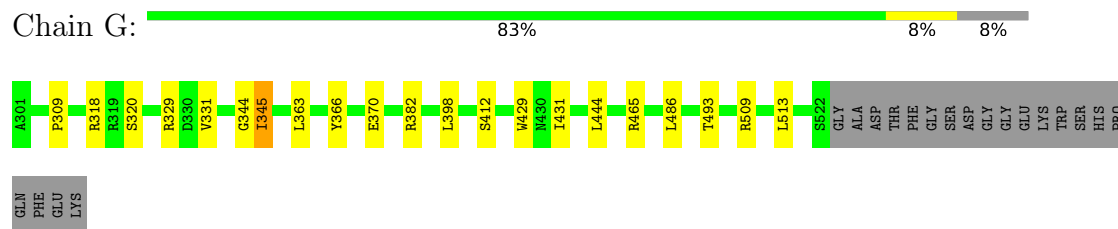
● Molecule 3: Proteasome subunit beta



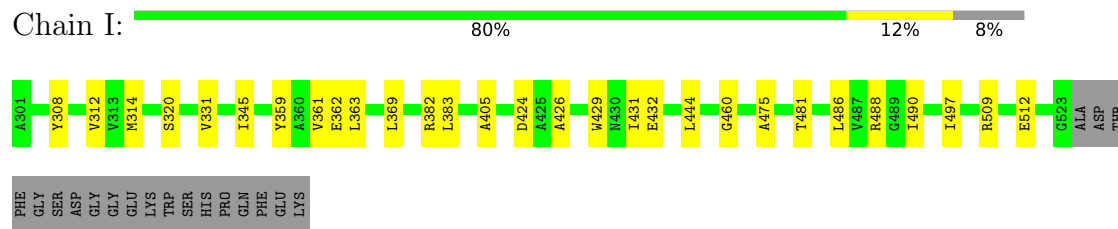
● Molecule 3: Proteasome subunit beta



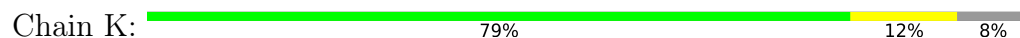
● Molecule 3: Proteasome subunit beta

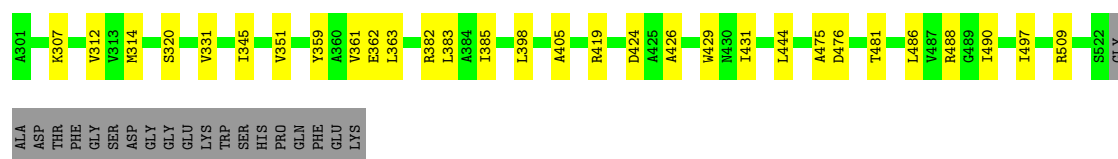


● Molecule 3: Proteasome subunit beta

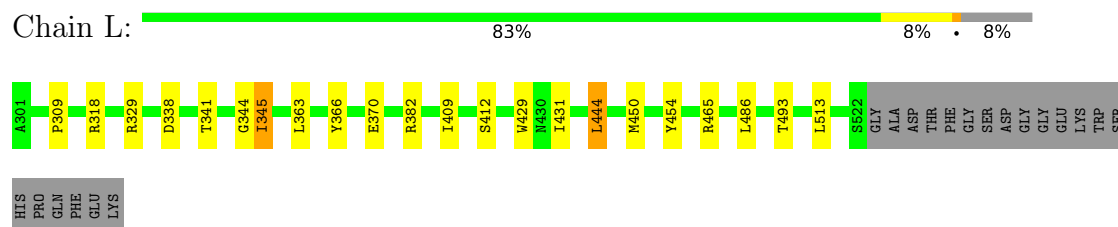


● Molecule 3: Proteasome subunit beta

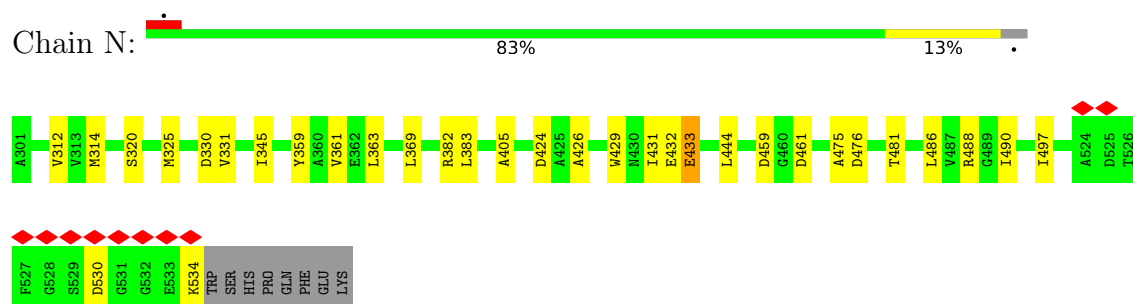




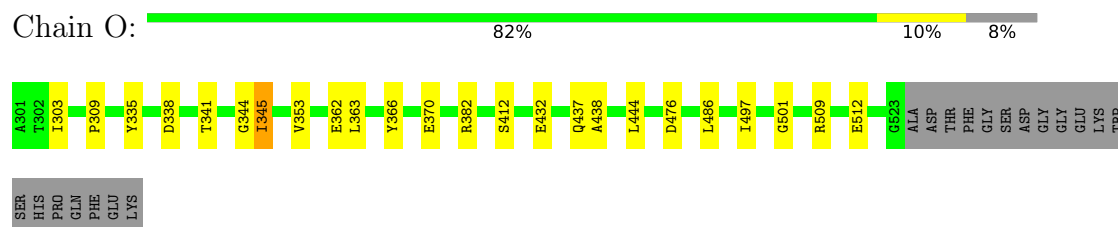
- Molecule 3: Proteasome subunit beta



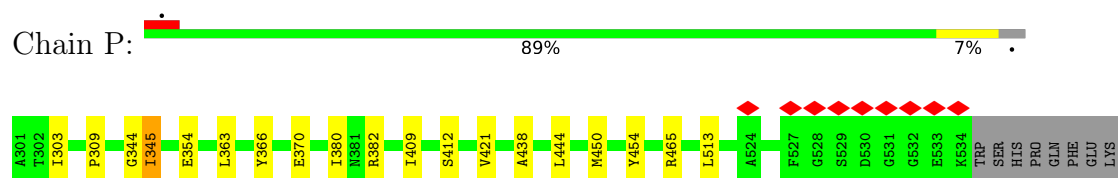
- Molecule 3: Proteasome subunit beta



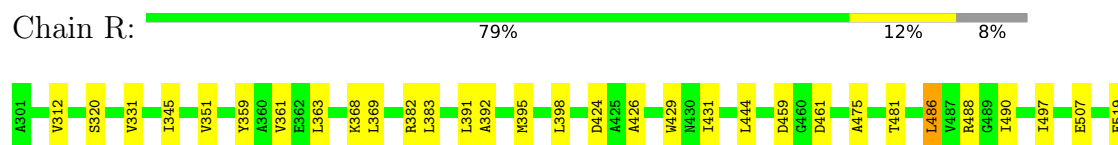
- Molecule 3: Proteasome subunit beta

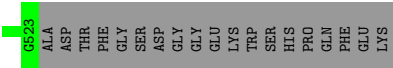


- Molecule 3: Proteasome subunit beta

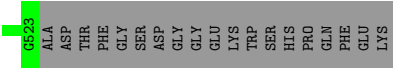
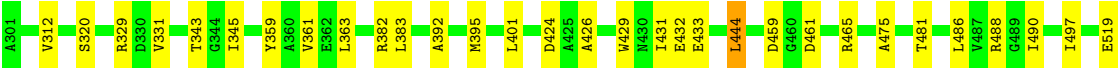
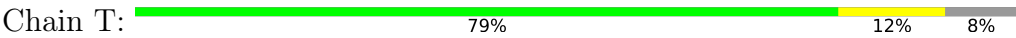


- Molecule 3: Proteasome subunit beta

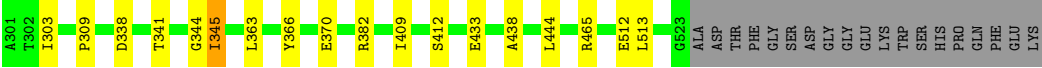
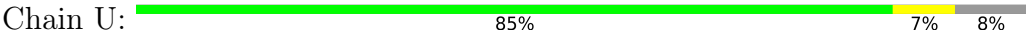




● Molecule 3: Proteasome subunit beta



● Molecule 3: Proteasome subunit beta



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C7	Depositor
Number of particles used	48799	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 ($e^-/\text{\AA}^2$)	25	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3600	Depositor
Magnification	100000	Depositor
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.395	Depositor
Minimum map value	-0.158	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.045	Depositor
Map size (Å)	537.6, 537.6, 537.6	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.4, 1.4, 1.4	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	0	0.34	0/1683	0.59	0/2274
1	2	0.34	0/1683	0.59	0/2274
1	4	0.33	0/1683	0.60	0/2274
1	6	0.33	0/1683	0.58	0/2274
1	8	0.33	0/1683	0.58	0/2274
1	B	0.32	0/1675	0.56	1/2263 (0.0%)
1	D	0.28	0/1693	0.52	0/2287
1	H	0.29	0/1683	0.54	0/2274
1	J	0.32	0/1683	0.53	0/2274
1	M	0.30	0/1675	0.52	0/2263
1	Q	0.30	0/1675	0.54	0/2263
1	S	0.38	1/1683 (0.1%)	0.55	1/2274 (0.0%)
1	W	0.33	0/1683	0.59	0/2274
1	Y	0.35	0/1683	0.59	0/2274
2	1	1.04	0/39	1.23	0/50
2	3	0.98	0/39	1.12	0/50
2	5	0.72	0/39	1.04	0/50
2	7	0.75	0/39	1.13	0/50
2	V	0.93	0/39	1.43	0/50
2	X	0.83	0/39	1.16	0/50
2	Z	1.04	0/39	1.28	0/50
3	A	0.33	0/1660	0.58	1/2251 (0.0%)
3	C	0.34	0/1664	0.58	2/2256 (0.1%)
3	E	0.37	1/1660 (0.1%)	0.56	0/2251
3	F	0.33	0/1660	0.56	1/2251 (0.0%)
3	G	0.34	0/1660	0.55	0/2251
3	I	0.34	0/1664	0.58	0/2256
3	K	0.35	0/1660	0.56	0/2251
3	L	0.33	0/1660	0.54	0/2251
3	N	0.36	0/1740	0.58	0/2357
3	O	0.35	0/1664	0.57	0/2256
3	P	0.33	0/1740	0.54	0/2357
3	R	0.35	0/1664	0.60	0/2256
3	T	0.36	0/1664	0.59	2/2256 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
3	U	0.33	0/1664	0.54	0/2256
All	All	0.34	2/47245 (0.0%)	0.57	8/63922 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	S	89	TYR	CA-C	-8.21	1.47	1.52
3	E	344	GLY	CA-C	-5.23	1.48	1.52

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	505	VAL	N-CA-C	-8.83	101.04	108.63
3	F	505	VAL	N-CA-C	-6.73	102.36	108.95
1	S	66	GLY	N-CA-C	5.50	118.64	110.60
1	B	66	GLY	N-CA-C	5.19	118.18	110.60
3	C	343	THR	CA-C-N	-5.14	117.81	122.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	0	1658	0	1659	34	0
1	2	1658	0	1659	33	0
1	4	1658	0	1659	34	0
1	6	1658	0	1659	29	0
1	8	1658	0	1659	34	0
1	B	1650	0	1648	36	0
1	D	1668	0	1667	28	0
1	H	1658	0	1659	24	0
1	J	1658	0	1659	29	0
1	M	1650	0	1648	33	0
1	Q	1650	0	1648	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	S	1658	0	1659	32	0
1	W	1658	0	1659	34	0
1	Y	1658	0	1659	35	0
2	1	39	0	32	5	0
2	3	39	0	32	5	0
2	5	39	0	32	5	0
2	7	39	0	32	5	0
2	V	39	0	32	7	0
2	X	39	0	32	5	0
2	Z	39	0	32	5	0
3	A	1636	0	1628	12	0
3	C	1640	0	1631	17	0
3	E	1636	0	1628	14	0
3	F	1636	0	1628	12	0
3	G	1636	0	1628	9	0
3	I	1640	0	1631	14	0
3	K	1636	0	1628	16	0
3	L	1636	0	1628	11	0
3	N	1715	0	1693	19	0
3	O	1640	0	1631	12	0
3	P	1715	0	1693	7	0
3	R	1640	0	1631	15	0
3	T	1640	0	1631	17	0
3	U	1640	0	1631	7	0
All	All	46557	0	46365	593	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 6.

The worst 5 of 593 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)
2:5:170:THR:N	1:8:13:MET:SD	2.46	0.89
2:3:170:THR:N	1:6:13:MET:SD	2.47	0.87
2:V:170:THR:N	1:Y:13:MET:SD	2.48	0.86
2:1:170:THR:N	1:4:13:MET:SD	2.48	0.86
2:7:170:THR:N	1:W:13:MET:SD	2.51	0.84

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	0	67/241 (28%)	65 (97%)	2 (3%)	0	100	100
1	2	211/241 (88%)	203 (96%)	8 (4%)	0	100	100
1	4	211/241 (88%)	203 (96%)	8 (4%)	0	100	100
1	6	151/241 (63%)	144 (95%)	7 (5%)	0	100	100
1	B	107/241 (44%)	102 (95%)	5 (5%)	0	100	100
1	D	213/241 (88%)	205 (96%)	8 (4%)	0	100	100
1	H	211/241 (88%)	205 (97%)	6 (3%)	0	100	100
1	J	211/241 (88%)	204 (97%)	7 (3%)	0	100	100
1	M	210/241 (87%)	202 (96%)	8 (4%)	0	100	100
1	Q	210/241 (87%)	204 (97%)	6 (3%)	0	100	100
1	S	211/241 (88%)	204 (97%)	7 (3%)	0	100	100
1	W	211/241 (88%)	201 (95%)	10 (5%)	0	100	100
1	Y	211/241 (88%)	201 (95%)	10 (5%)	0	100	100
2	1	3/180 (2%)	3 (100%)	0	0	100	100
2	3	3/180 (2%)	3 (100%)	0	0	100	100
2	5	3/180 (2%)	3 (100%)	0	0	100	100
2	V	3/180 (2%)	3 (100%)	0	0	100	100
2	X	3/180 (2%)	3 (100%)	0	0	100	100
2	Z	3/180 (2%)	3 (100%)	0	0	100	100
3	C	221/242 (91%)	211 (96%)	10 (4%)	0	100	100
3	E	220/242 (91%)	209 (95%)	11 (5%)	0	100	100
3	F	220/242 (91%)	211 (96%)	8 (4%)	1 (0%)	24	57
3	G	220/242 (91%)	212 (96%)	7 (3%)	1 (0%)	24	57
3	I	221/242 (91%)	210 (95%)	11 (5%)	0	100	100
3	K	220/242 (91%)	210 (96%)	10 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	L	220/242 (91%)	213 (97%)	6 (3%)	1 (0%)	24	57
3	N	232/242 (96%)	221 (95%)	11 (5%)	0	100	100
3	O	221/242 (91%)	214 (97%)	6 (3%)	1 (0%)	24	57
3	P	232/242 (96%)	225 (97%)	6 (3%)	1 (0%)	30	62
3	R	221/242 (91%)	210 (95%)	11 (5%)	0	100	100
3	T	221/242 (91%)	210 (95%)	11 (5%)	0	100	100
3	U	221/242 (91%)	214 (97%)	6 (3%)	1 (0%)	24	57
All	All	5343/7359 (73%)	5131 (96%)	206 (4%)	6 (0%)	49	79

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	P	309	PRO
3	U	309	PRO
3	F	309	PRO
3	G	309	PRO
3	L	309	PRO

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	0	165/185 (89%)	154 (93%)	11 (7%)	15	41
1	2	165/185 (89%)	154 (93%)	11 (7%)	15	41
1	4	165/185 (89%)	154 (93%)	11 (7%)	15	41
1	6	165/185 (89%)	155 (94%)	10 (6%)	17	43
1	8	165/185 (89%)	154 (93%)	11 (7%)	15	41
1	B	164/185 (89%)	156 (95%)	8 (5%)	22	48
1	D	166/185 (90%)	158 (95%)	8 (5%)	23	49
1	H	165/185 (89%)	155 (94%)	10 (6%)	17	43
1	J	165/185 (89%)	153 (93%)	12 (7%)	13	38

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	M	164/185 (89%)	157 (96%)	7 (4%)	26	51
1	Q	164/185 (89%)	159 (97%)	5 (3%)	36	60
1	S	165/185 (89%)	157 (95%)	8 (5%)	23	49
1	W	165/185 (89%)	154 (93%)	11 (7%)	15	41
1	Y	165/185 (89%)	154 (93%)	11 (7%)	15	41
2	1	3/147 (2%)	3 (100%)	0	100	100
2	3	3/147 (2%)	3 (100%)	0	100	100
2	5	3/147 (2%)	3 (100%)	0	100	100
2	7	3/147 (2%)	3 (100%)	0	100	100
2	V	3/147 (2%)	3 (100%)	0	100	100
2	X	3/147 (2%)	3 (100%)	0	100	100
2	Z	3/147 (2%)	3 (100%)	0	100	100
3	A	164/179 (92%)	159 (97%)	5 (3%)	36	60
3	C	164/179 (92%)	157 (96%)	7 (4%)	26	51
3	E	164/179 (92%)	158 (96%)	6 (4%)	30	56
3	F	164/179 (92%)	157 (96%)	7 (4%)	26	51
3	G	164/179 (92%)	158 (96%)	6 (4%)	30	56
3	I	164/179 (92%)	158 (96%)	6 (4%)	30	56
3	K	164/179 (92%)	159 (97%)	5 (3%)	36	60
3	L	164/179 (92%)	158 (96%)	6 (4%)	30	56
3	N	171/179 (96%)	164 (96%)	7 (4%)	27	53
3	O	164/179 (92%)	157 (96%)	7 (4%)	26	51
3	P	171/179 (96%)	165 (96%)	6 (4%)	32	57
3	R	164/179 (92%)	157 (96%)	7 (4%)	26	51
3	T	164/179 (92%)	158 (96%)	6 (4%)	30	56
3	U	164/179 (92%)	157 (96%)	7 (4%)	26	51
All	All	4639/6125 (76%)	4417 (95%)	222 (5%)	24	49

5 of 222 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	I	444	LEU
1	Y	234	LEU

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Mol	Chain	Res	Type
1	M	74	LEU
1	Y	203	LEU
3	U	512	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 95 such sidechains are listed below:

Mol	Chain	Res	Type
1	J	114	GLN
1	Q	101	ASN
3	K	456	GLN
3	O	456	GLN
1	S	69	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.

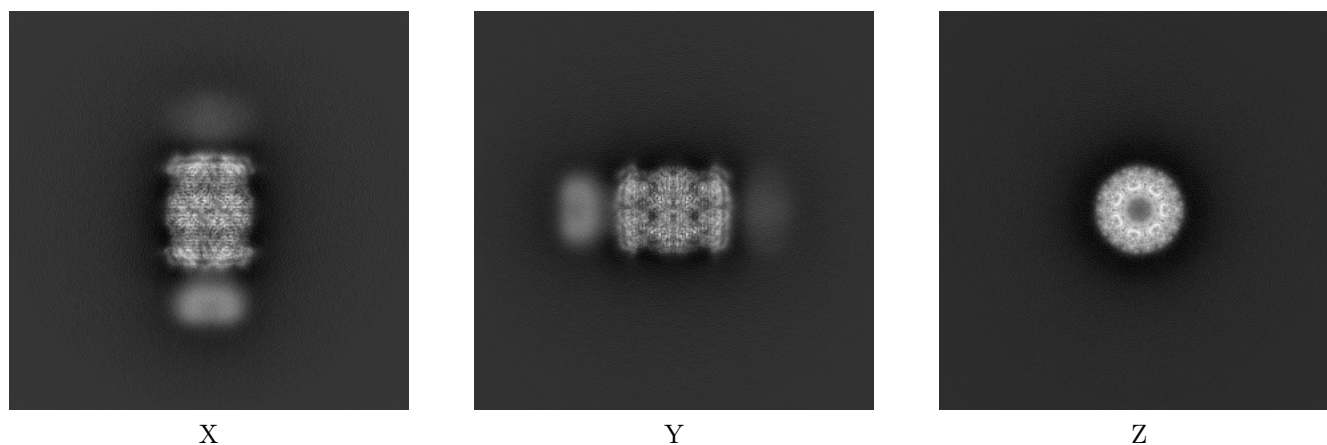
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-4128. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

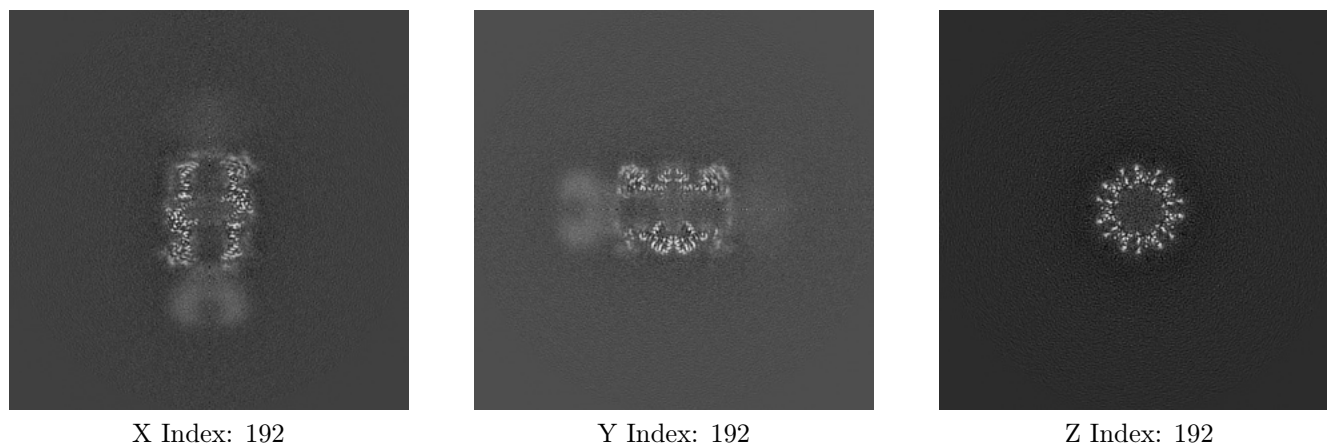
6.1.1 Primary map



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

6.2 Central slices [i](#)

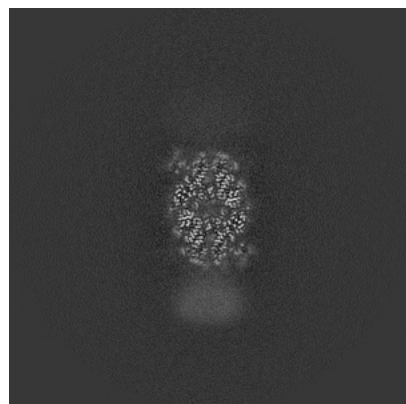
6.2.1 Primary map



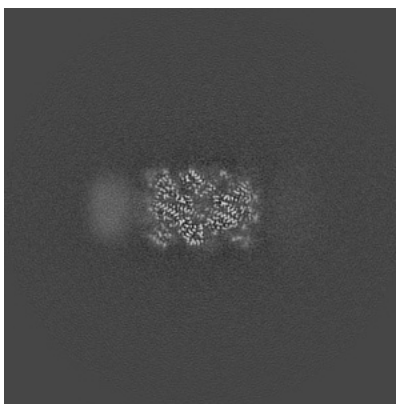
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

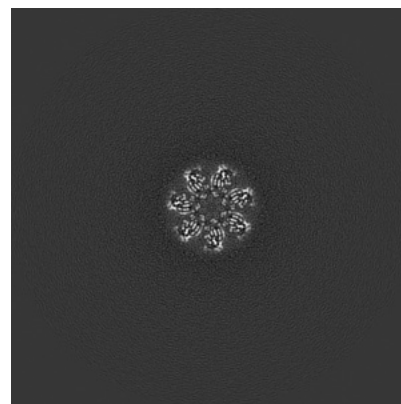
6.3.1 Primary map



X Index: 174



Y Index: 171

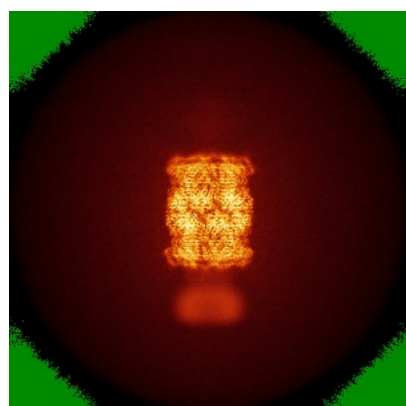


Z Index: 178

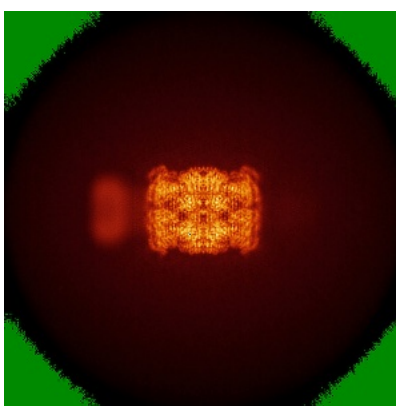
The images above show the largest variance slices of the map in three orthogonal directions.

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

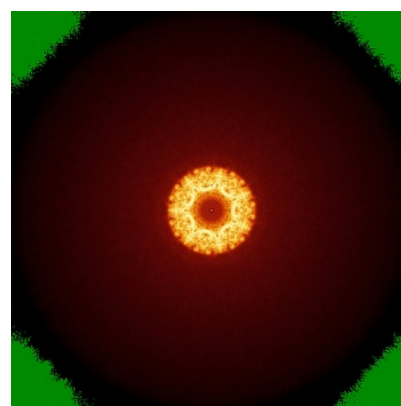
6.4.1 Primary map



X



Y

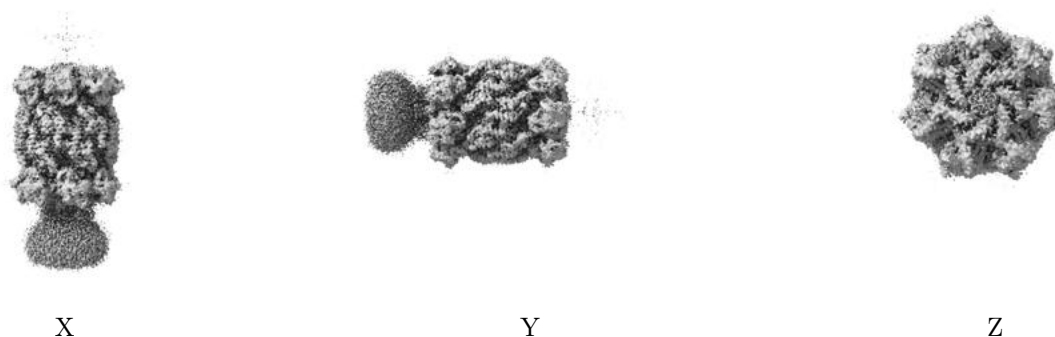


Z

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 0.045. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

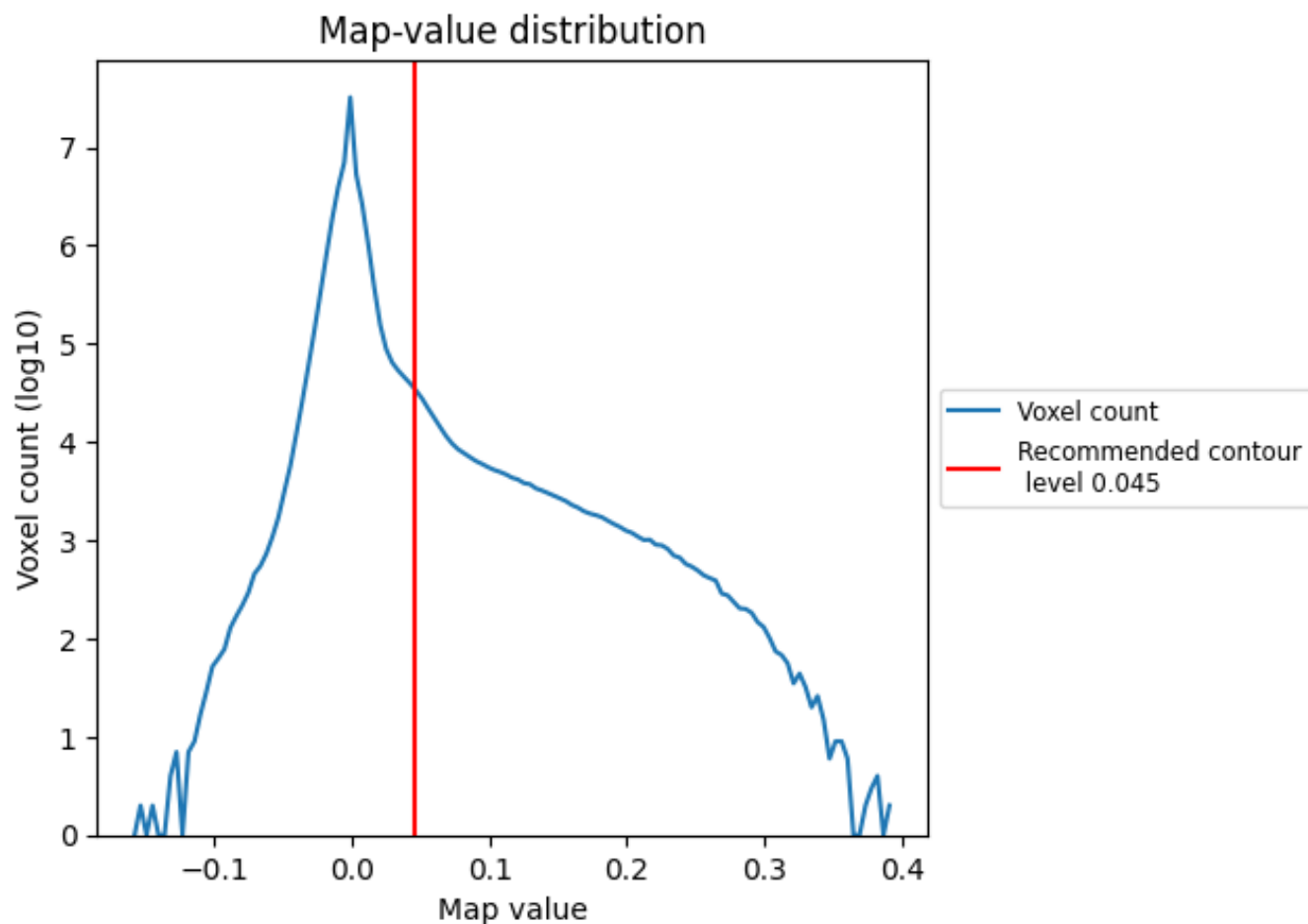
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

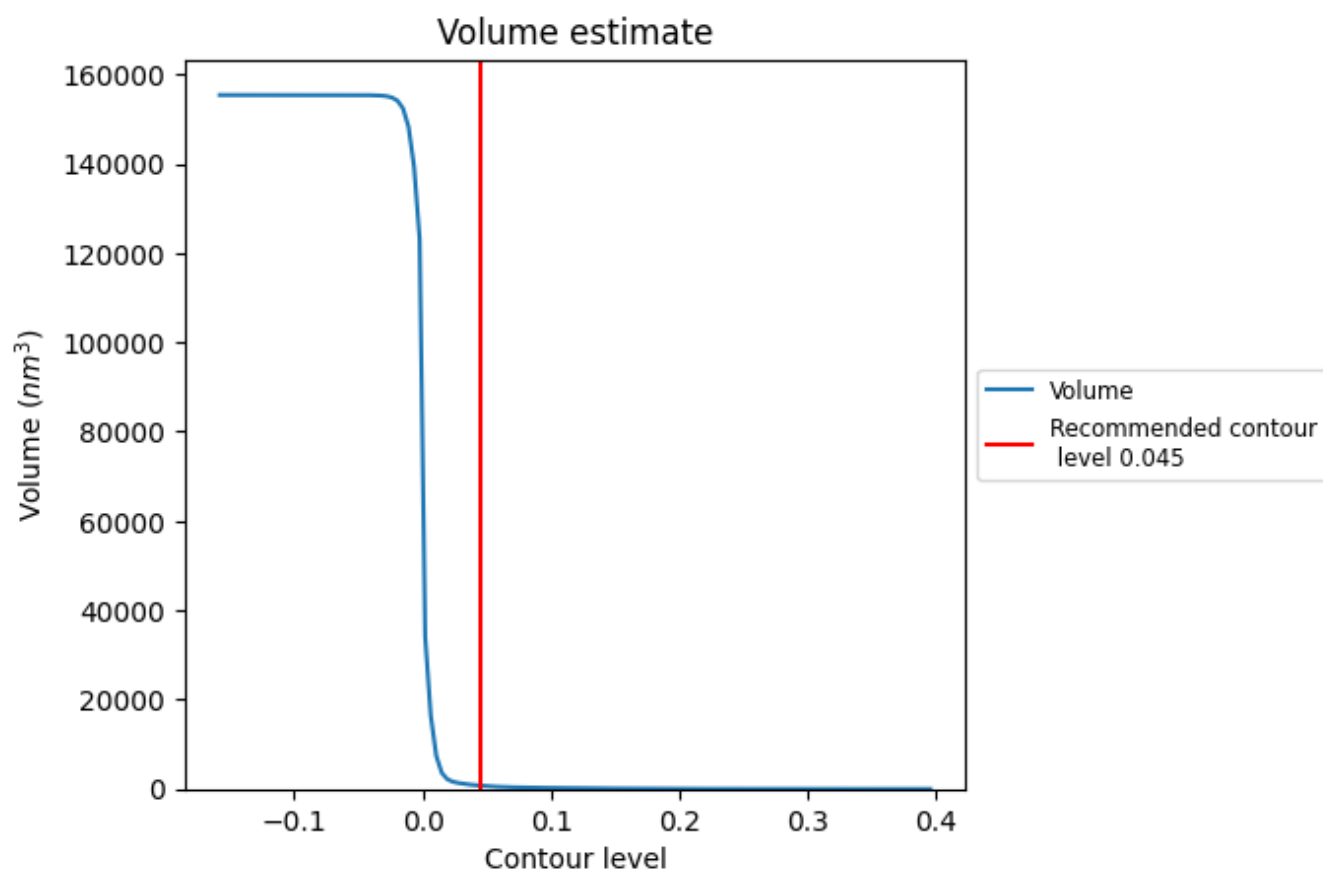
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

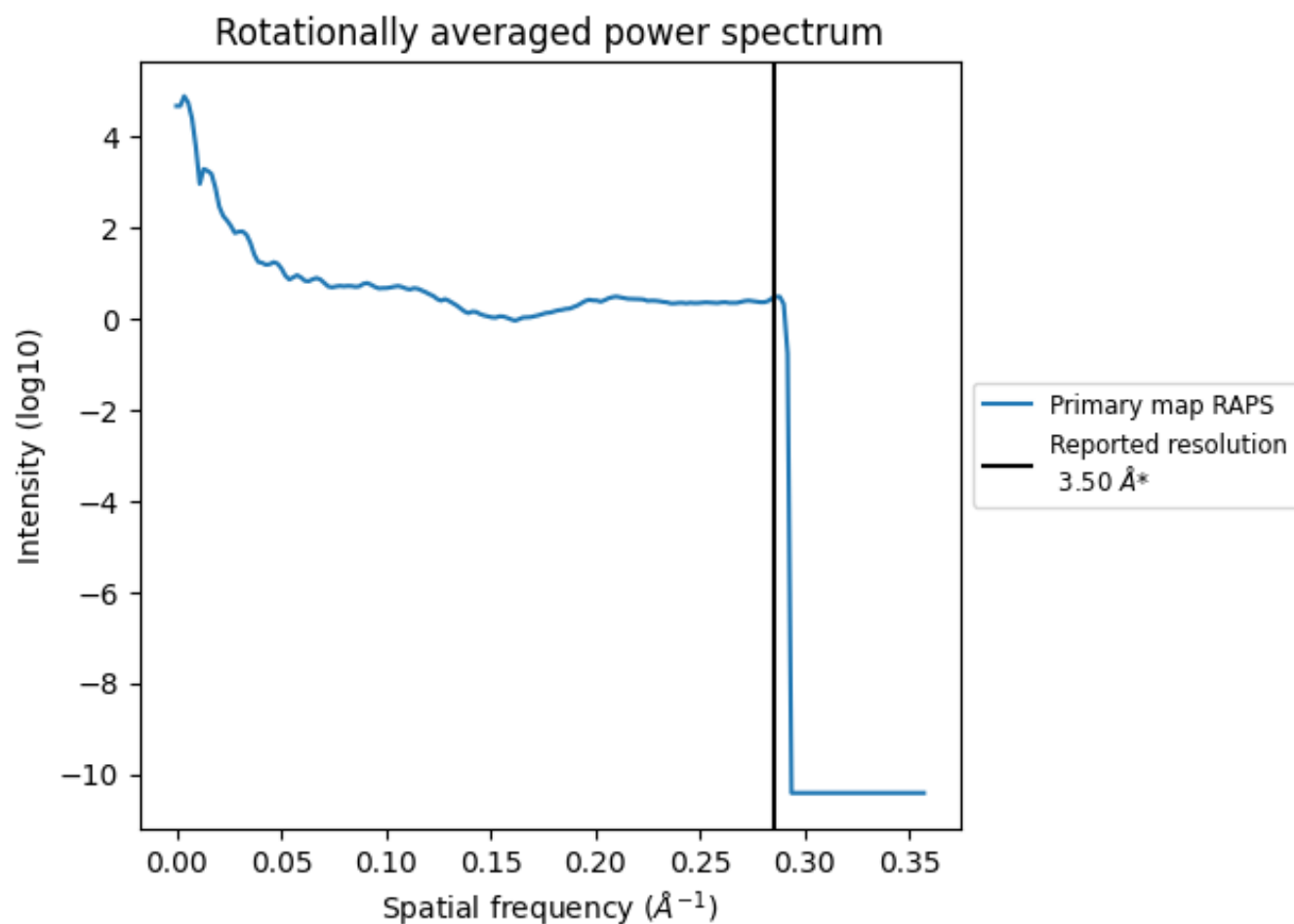
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 746 nm³; this corresponds to an approximate mass of 674 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 ⓘ

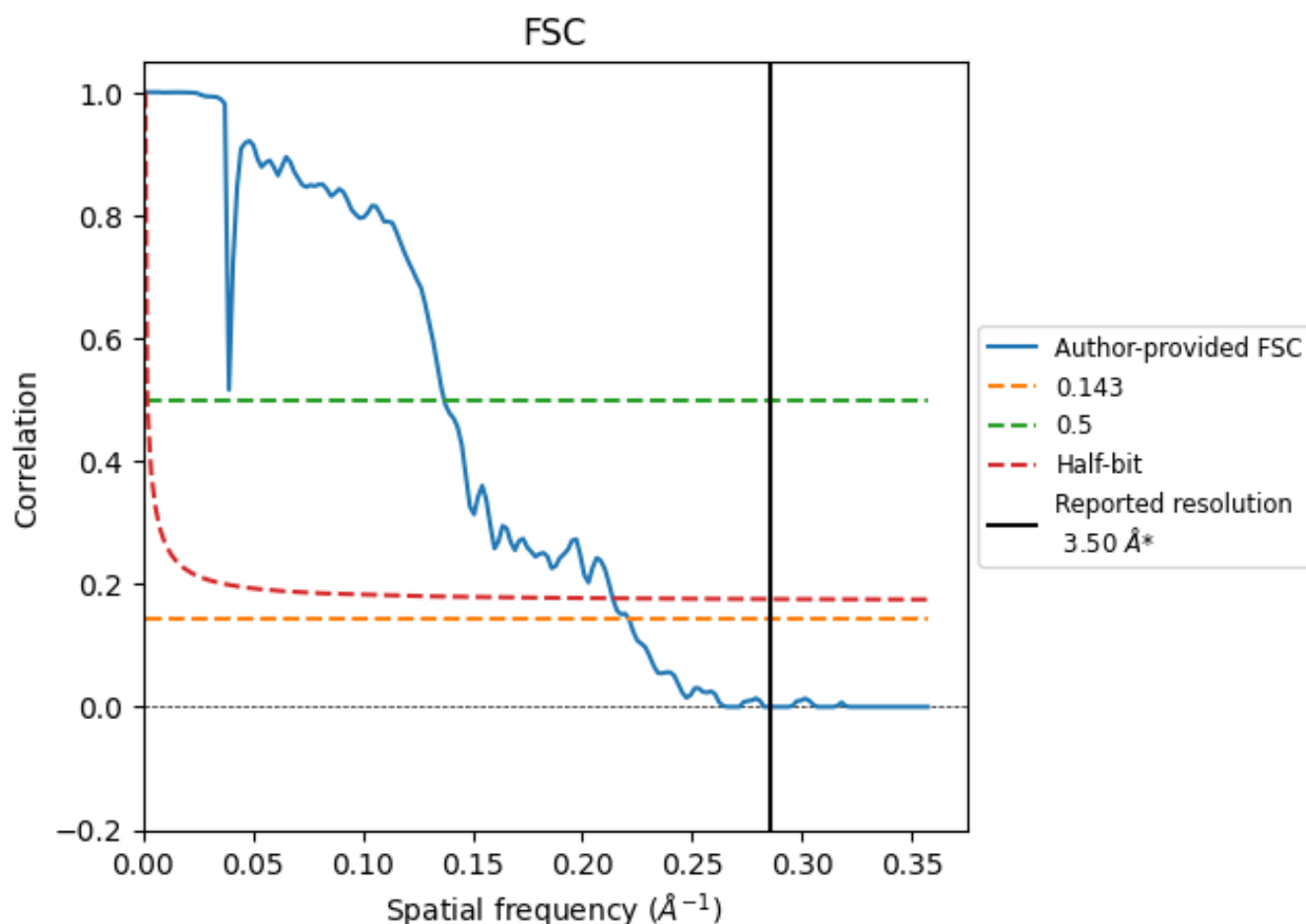


*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

8.2 Resolution estimates [i](#)

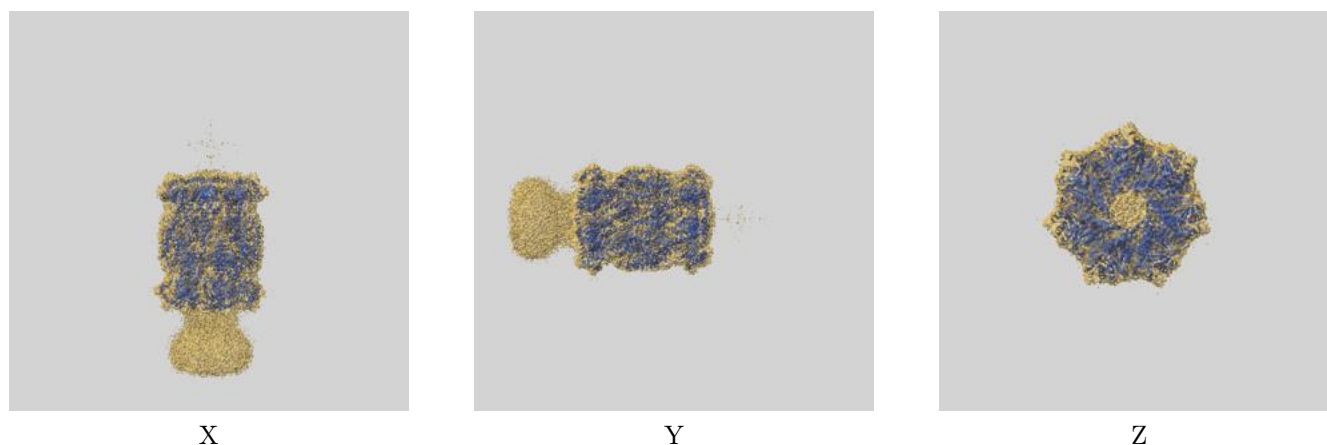
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	4.52	7.29	4.68
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from author-provided FSC intersecting FSC 0.143 CUT-OFF 4.52 differs from the reported value 3.5 by more than 10 %

9 Map-model fit [i](#)

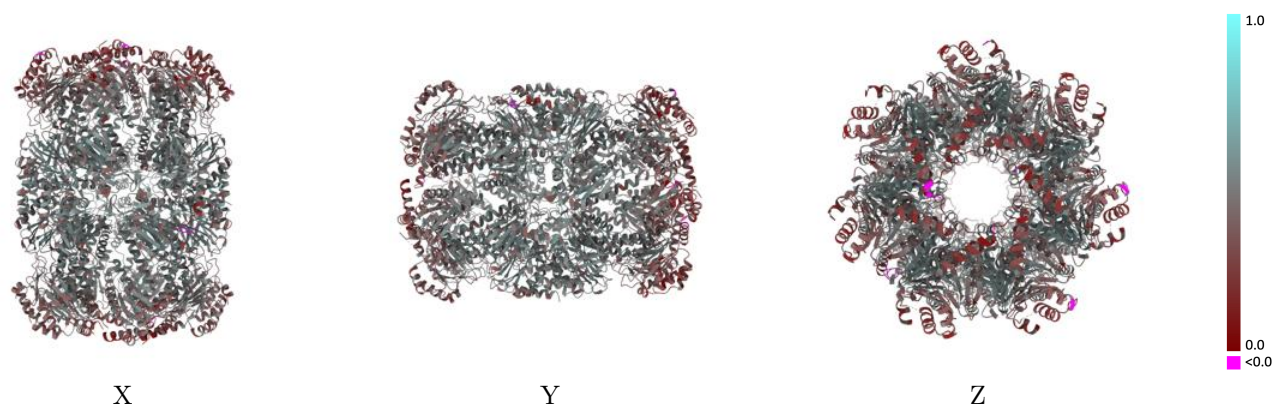
This section contains information regarding the fit between EMDB map EMD-4128 and PDB model 5LZP. Per-residue inclusion information can be found in section [3](#) on page [12](#).

9.1 Map-model overlay [i](#)



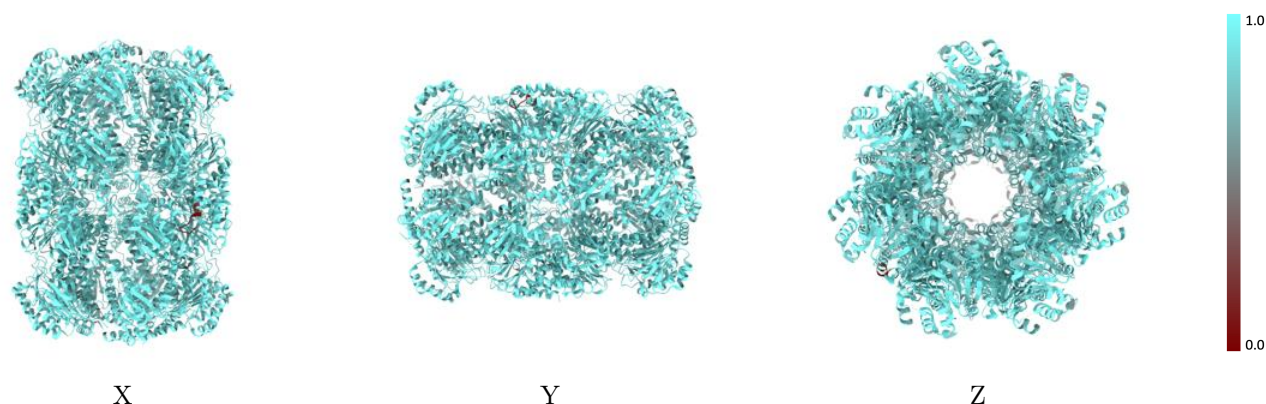
The images above show the 3D surface view of the map at the recommended contour level 0.045 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



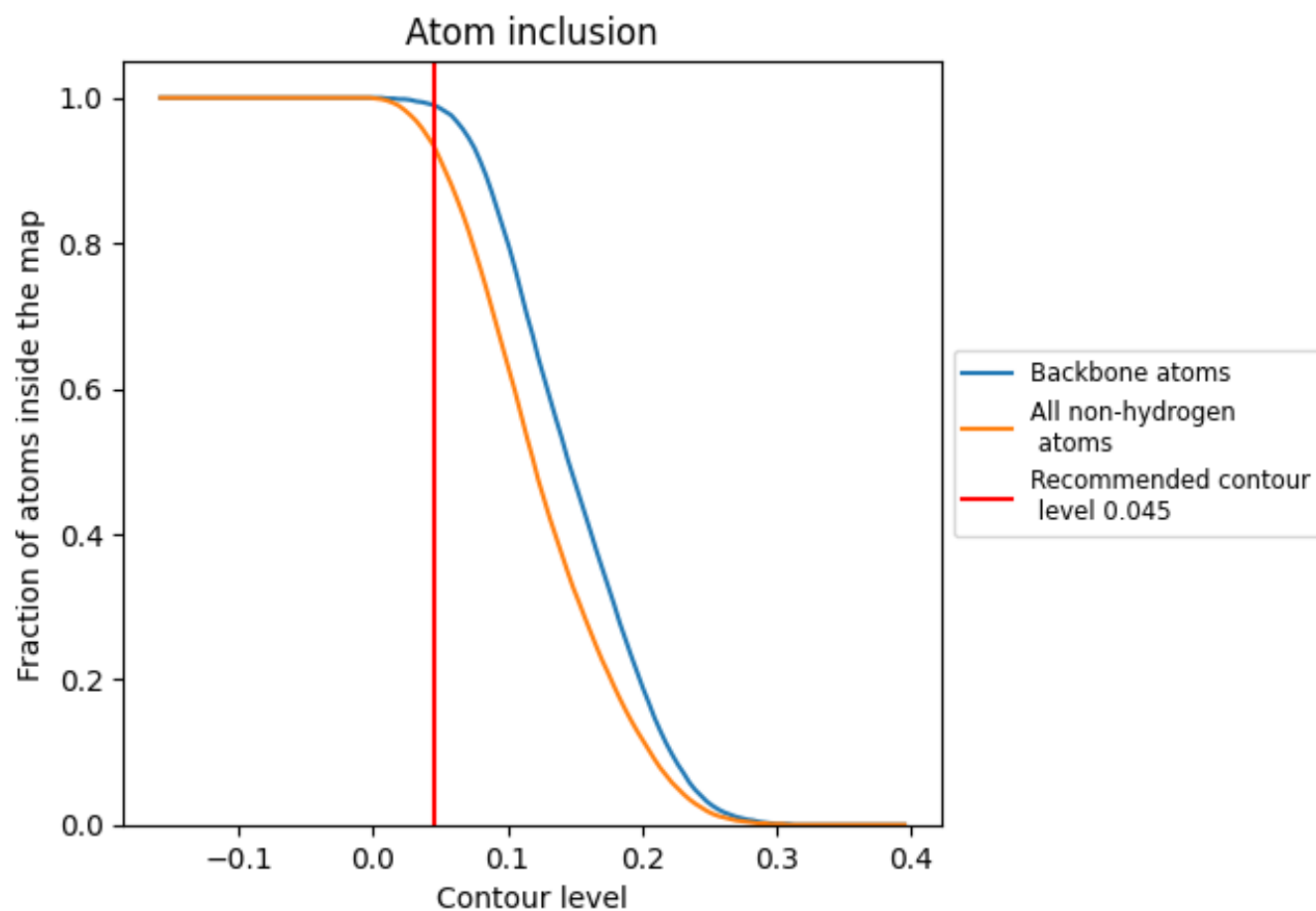
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.045).

9.4 Atom inclusion ⓘ



At the recommended contour level, 99% of all backbone atoms, 93% 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 (0.045) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9340	 0.4400
0	 0.9210	 0.4130
1	 0.8160	 0.2610
2	 0.9230	 0.4130
3	 0.7890	 0.2580
4	 0.9250	 0.4120
5	 0.7630	 0.2420
6	 0.9260	 0.4120
7	 0.7890	 0.2550
8	 0.9200	 0.4130
A	 0.9570	 0.4930
B	 0.9230	 0.3700
C	 0.9450	 0.4910
D	 0.9310	 0.3760
E	 0.9530	 0.4880
F	 0.9560	 0.4930
G	 0.9560	 0.4930
H	 0.9270	 0.3760
I	 0.9460	 0.4890
J	 0.9250	 0.3720
K	 0.9480	 0.4910
L	 0.9560	 0.4960
M	 0.9220	 0.3780
N	 0.9100	 0.4710
O	 0.9540	 0.4930
P	 0.9180	 0.4690
Q	 0.9230	 0.3650
R	 0.9480	 0.4880
S	 0.9250	 0.3740
T	 0.9460	 0.4910
U	 0.9550	 0.4930
V	 0.7890	 0.2300
W	 0.9210	 0.4150
X	 0.8160	 0.2460
Y	 0.9260	 0.4140
Z	 0.7890	 0.2280

