



## wwPDB EM Validation Summary Report ⓘ

Oct 31, 2022 – 04:35 PM EDT

PDB ID : 8D6Y  
EMDB ID : EMD-27226  
Title : Structure of the Mycobacterium tuberculosis 20S proteasome bound to the ADP-bound Mpa ATPase  
Authors : Xiao, X.; Li, H.  
Deposited on : 2022-06-06  
Resolution : 10.00 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

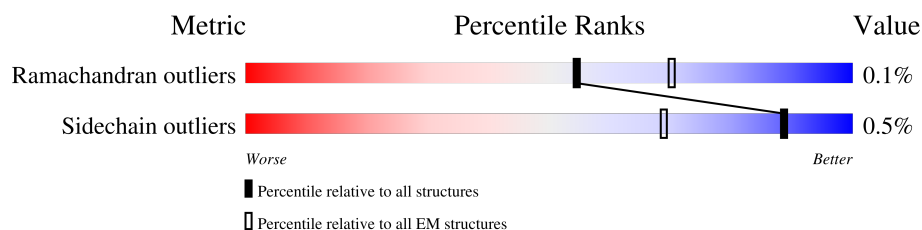
EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

# 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 10.00 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.









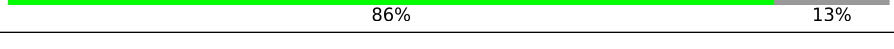
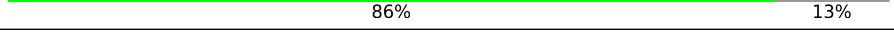
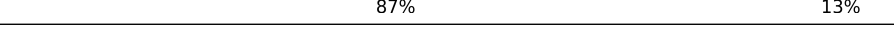
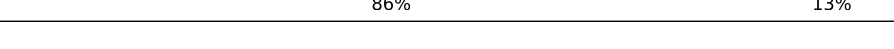
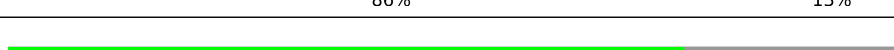

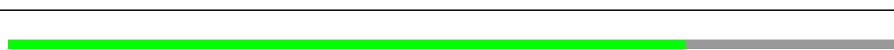

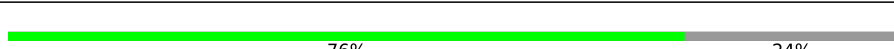





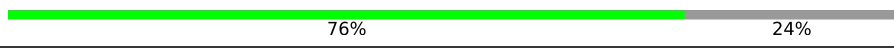
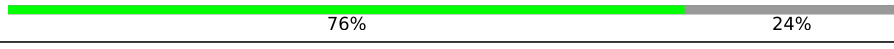



Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	609	<div> <div>13%</div> <div>78%</div> <div>21%</div> </div>
1	B	609	<div> <div>11%</div> <div>77%</div> <div>22%</div> </div>
1	C	609	<div> <div>8%</div> <div>77%</div> <div>22%</div> </div>
1	D	609	<div> <div>10%</div> <div>75%</div> <div>24%</div> </div>
1	E	609	<div> <div>13%</div> <div>75%</div> <div>24%</div> </div>
1	F	609	<div> <div>16%</div> <div>75%</div> <div>24%</div> </div>
2	G	248	<div> <div>86%</div> <div>13%</div> </div>
2	H	248	<div> <div>86%</div> <div>13%</div> </div>
2	I	248	<div> <div>86%</div> <div>13%</div> </div>

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Mol	Chain	Length	Quality of chain
2	J	248	
2	K	248	
2	L	248	
2	M	248	
2	N	248	
2	O	248	
2	k	248	
2	l	248	
2	m	248	
2	n	248	
2	o	248	
3	P	291	
3	Q	291	
3	R	291	
3	S	291	
3	T	291	
3	U	291	
3	V	291	
3	W	291	
3	X	291	
3	Y	291	
3	Z	291	
3	a	291	
3	b	291	
3	c	291	

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Mol	Chain	Length	Quality of chain
4	d	4	<div><div>50%</div><div>100%</div></div>
4	e	4	<div><div>75%</div><div>100%</div></div>
4	f	4	<div><div>75%</div><div>100%</div></div>
4	g	4	<div><div>25%</div><div>100%</div></div>
4	h	4	<div><div>50%</div><div>100%</div></div>
4	i	4	<div><div>50%</div><div>100%</div></div>
4	j	4	<div><div>25%</div><div>100%</div></div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 68262 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 AAA ATPase forming ring-shaped complexes.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	480	Total	C	N	O	S	0	0
			3747	2360	647	729	11		
1	B	475	Total	C	N	O	S	0	0
			3707	2337	637	722	11		
1	C	472	Total	C	N	O	S	0	0
			3649	2300	632	706	11		
1	D	460	Total	C	N	O	S	0	0
			3593	2266	621	695	11		
1	E	461	Total	C	N	O	S	0	0
			3597	2268	622	696	11		
1	F	460	Total	C	N	O	S	0	0
			3589	2267	615	696	11		

- Molecule 2 is a protein called Proteasome subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	k	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	l	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	m	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	n	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	o	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	G	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	H	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	I	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	J	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	K	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	L	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	M	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	N	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		
2	O	215	Total	C	N	O	S	0	0
			1658	1039	303	313	3		

- Molecule 3 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	P	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	Q	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	R	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	S	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	T	222	Total	C	N	O	S	0	0
			1636	1026	282	323	5		
3	U	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	V	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	W	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	X	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	Y	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	Z	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	a	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	b	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		
3	c	222	Total	C	N	O	S	0	0
			1638	1027	282	324	5		

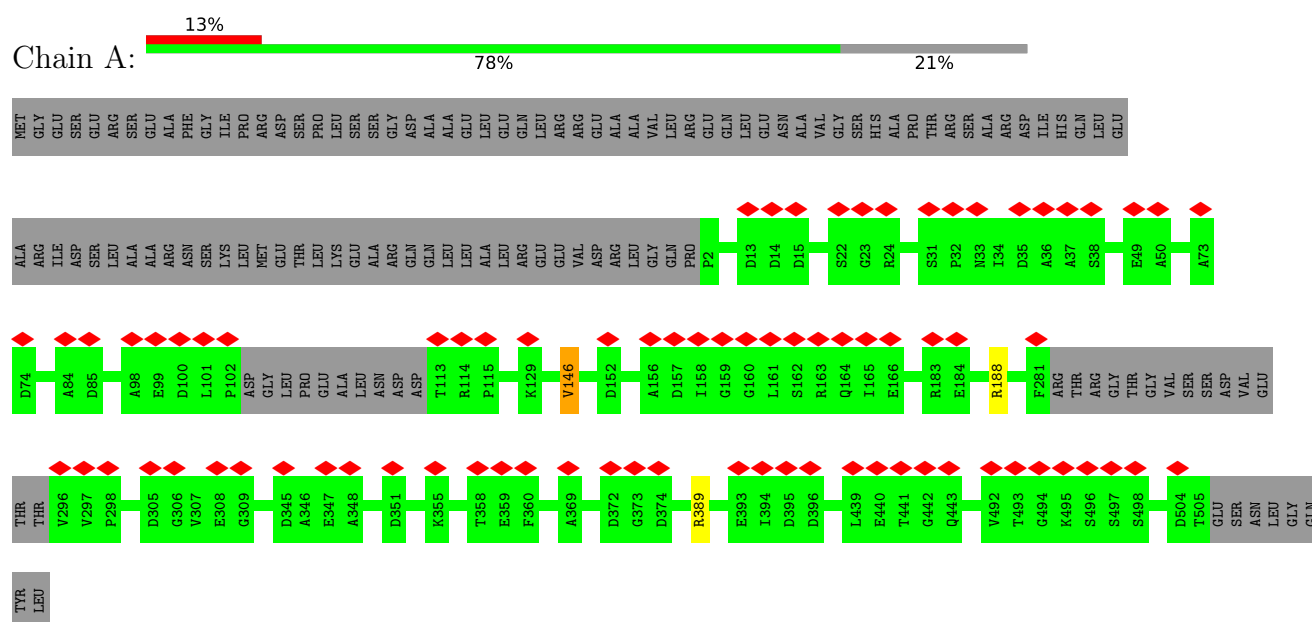
- Molecule 4 is a protein called Proteasome-associated ATPase.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	d	4	Total	C	N	O	0	0
			34	22	5	7		
4	e	4	Total	C	N	O	0	0
			34	22	5	7		
4	f	4	Total	C	N	O	0	0
			34	22	5	7		
4	g	4	Total	C	N	O	0	0
			34	22	5	7		
4	h	4	Total	C	N	O	0	0
			34	22	5	7		
4	i	4	Total	C	N	O	0	0
			34	22	5	7		
4	j	4	Total	C	N	O	0	0
			34	22	5	7		

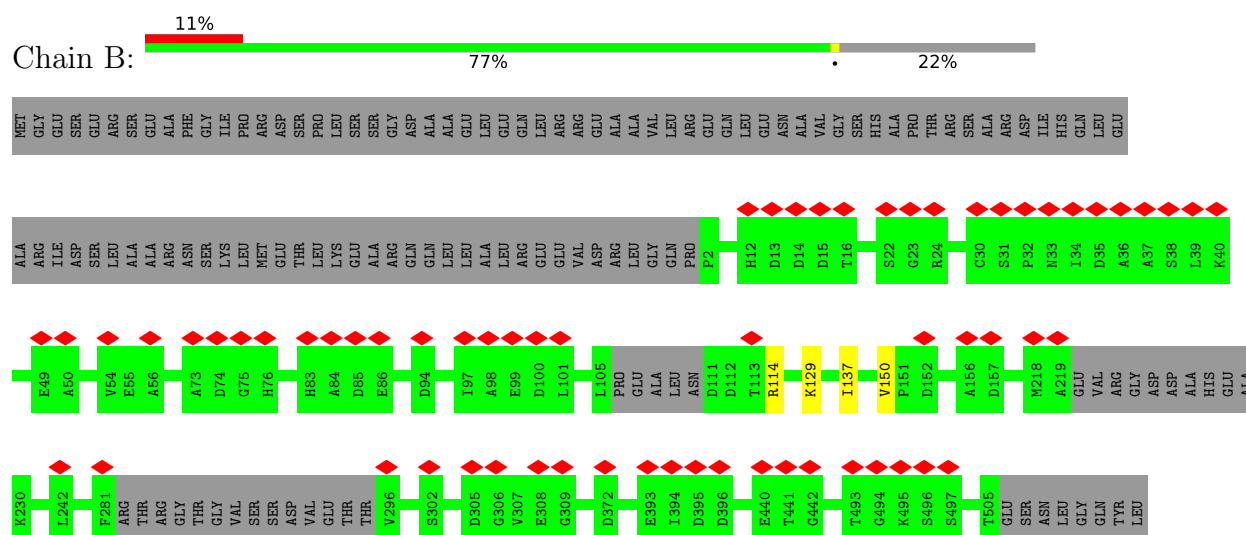
### 3 Residue-property plots

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: AAA ATPase forming ring-shaped complexes

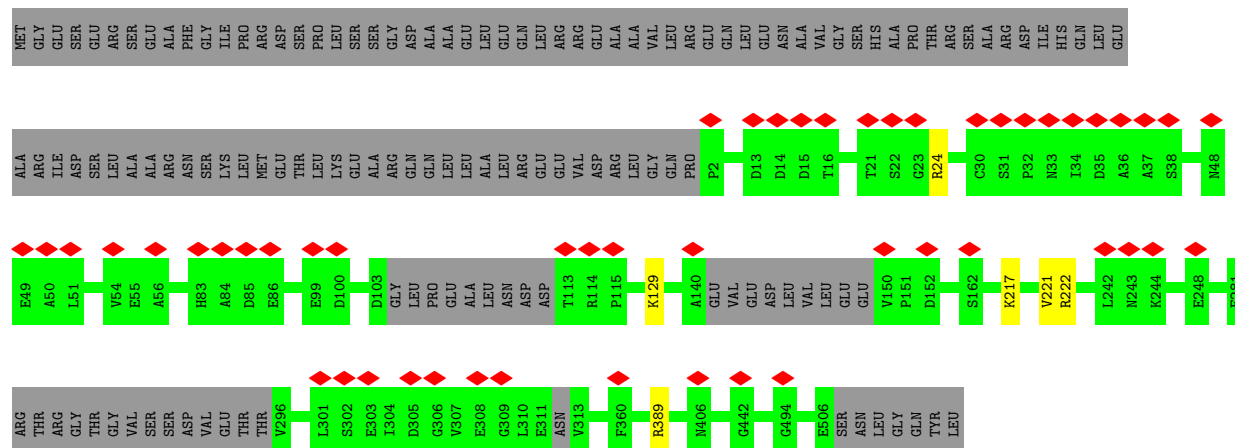
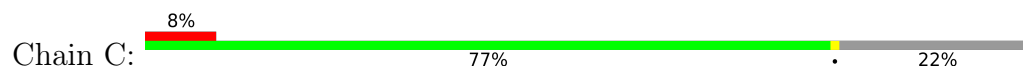


- Molecule 1: AAA ATPase forming ring-shaped complexes

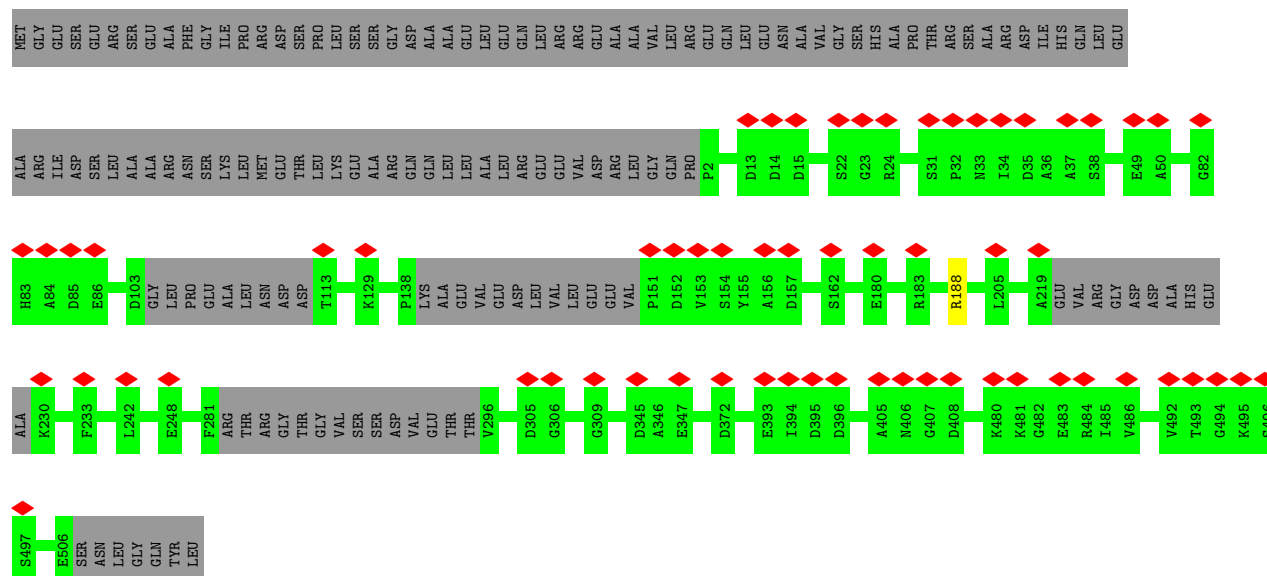
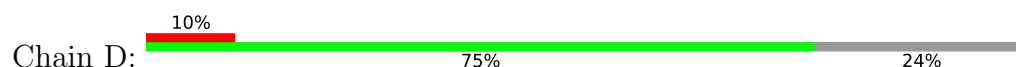




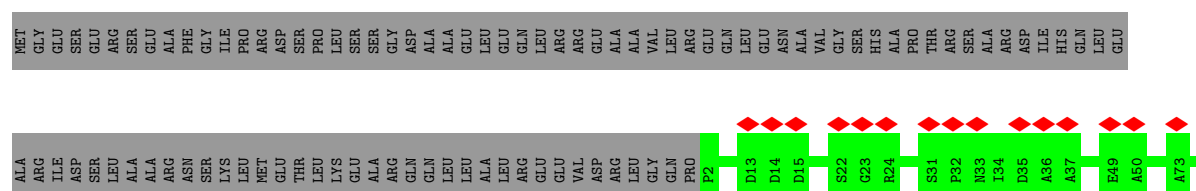
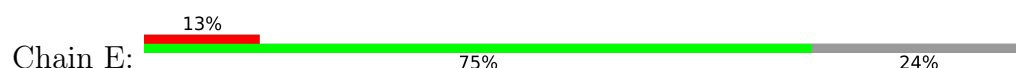
- Molecule 1: AAA ATPase forming ring-shaped complexes

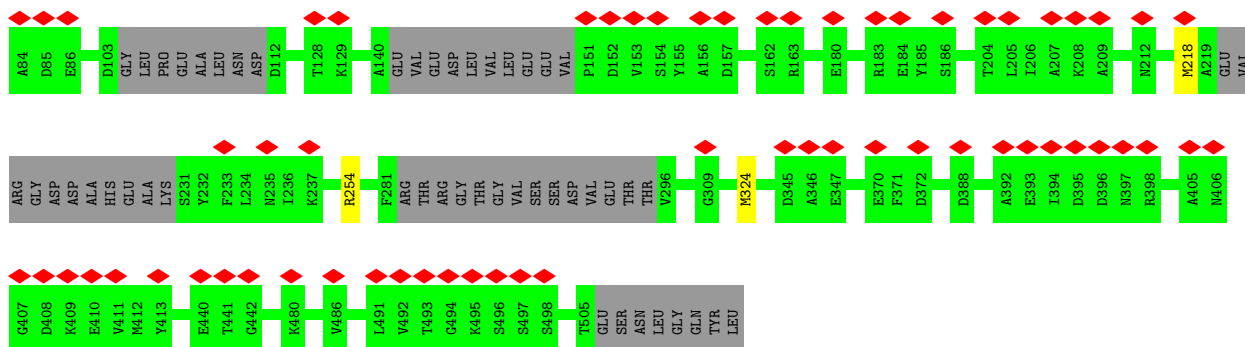


- Molecule 1: AAA ATPase forming ring-shaped complexes

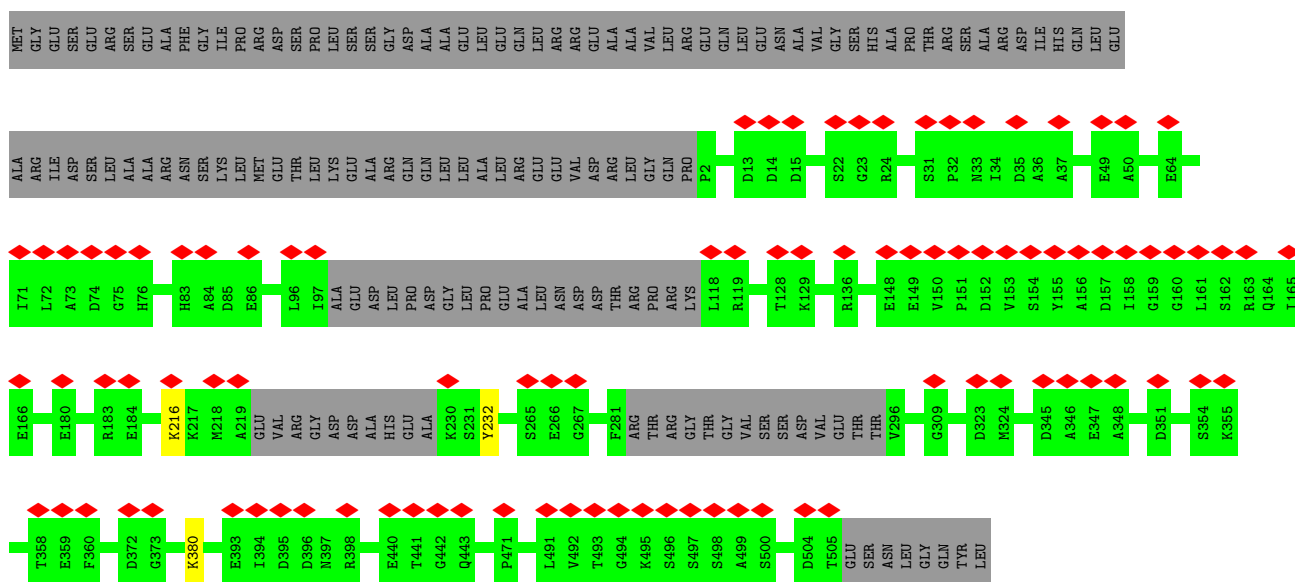
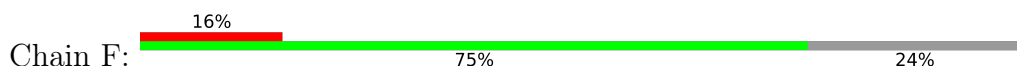


- Molecule 1: AAA ATPase forming ring-shaped complexes

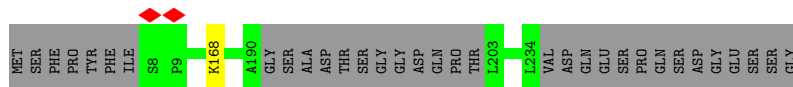
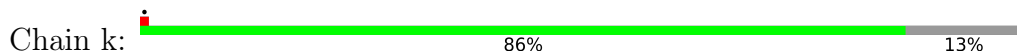




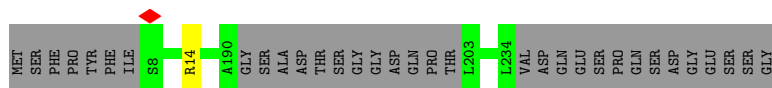
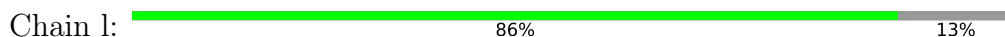
- Molecule 1: AAA ATPase forming ring-shaped complexes



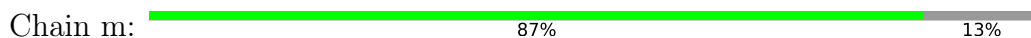
- Molecule 2: Proteasome subunit alpha

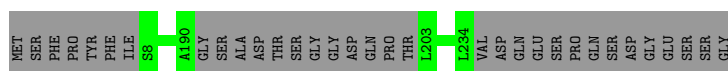


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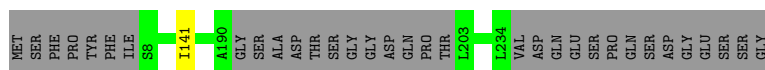
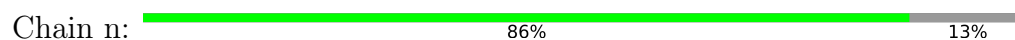


- Molecule 2: Proteasome subunit alpha

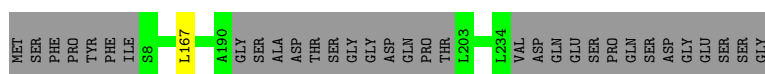
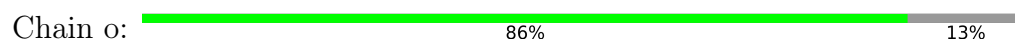




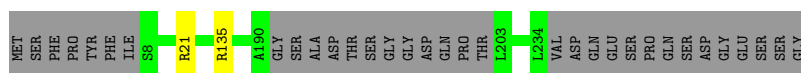
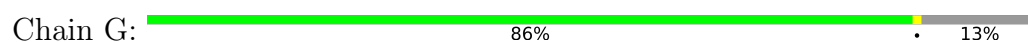
- Molecule 2: Proteasome subunit alpha



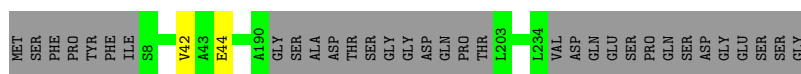
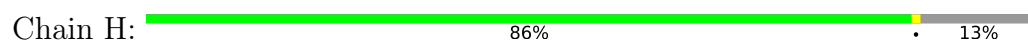
- Molecule 2: Proteasome subunit alpha



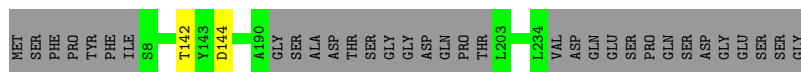
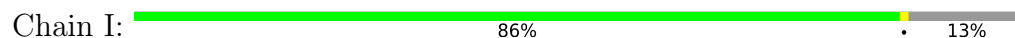
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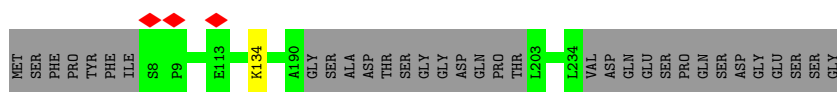
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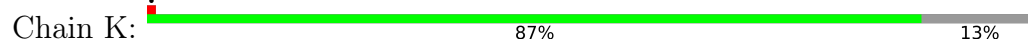
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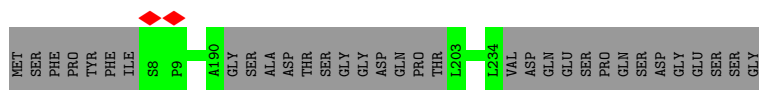


- Molecule 2: Proteasome subunit alpha

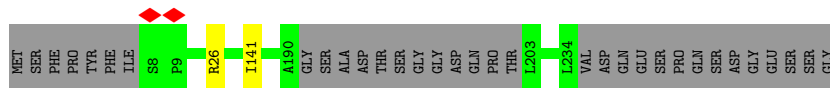
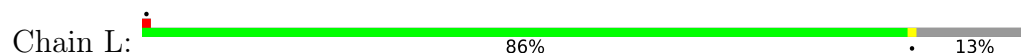


- Molecule 2: Proteasome subunit alpha

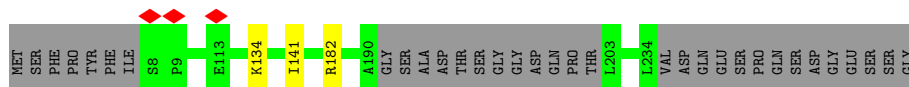
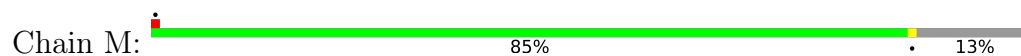




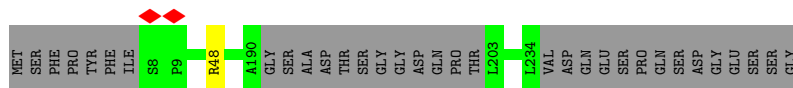
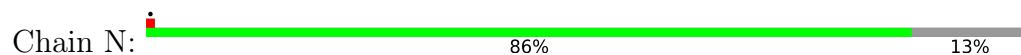
- Molecule 2: Proteasome subunit alpha



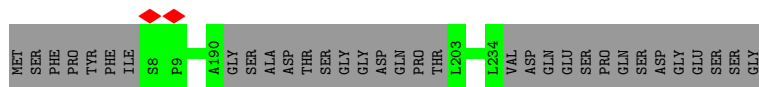
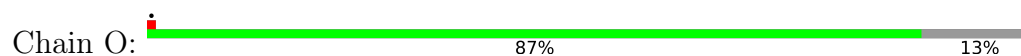
- Molecule 2: Proteasome subunit alpha



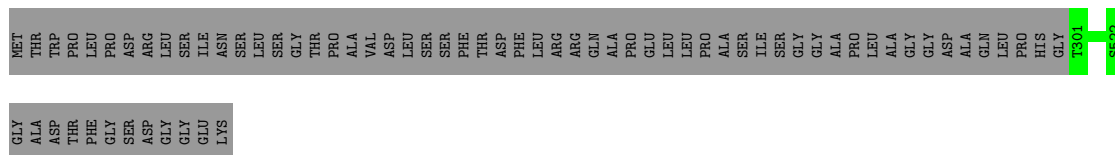
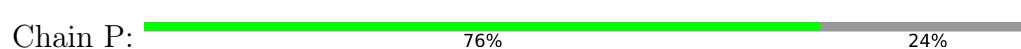
- Molecule 2: Proteasome subunit alpha



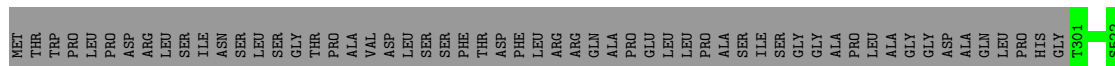
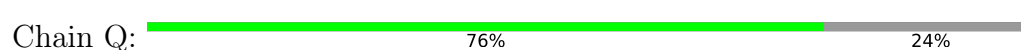
- Molecule 2: Proteasome subunit alpha



- Molecule 3: Proteasome subunit beta



- Molecule 3: Proteasome subunit beta



GLY  
ALA  
ASP  
THR  
PHE  
GLY  
SER  
ASP  
GLY  
GLU  
LYS

● Molecule 3: Proteasome subunit beta



MET THR TRP PRO LEU LEU PRO ASP ARG LEU SER ILE ASN SER LEU SER GLY THR PRO PRO ALA VAL ASP LEU SER SER PHE THR ASP PHE LEU ARG ARG GLN ALA PRO GLU LEU LEU PRO ALA ALA ILE SER SER GLY GLY ALA PRO LEU LEU GLY GLY ASP ALA GLN LEU LEU PRO HIS GLY T301 S522

GLY  
ALA  
ASP  
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GLY  
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GLY  
GLU  
LYS

● Molecule 3: Proteasome subunit beta



MET THR TRP PRO LEU LEU PRO ASP ARG LEU SER ILE ASN SER LEU SER GLY THR PRO PRO ALA VAL ASP LEU SER SER PHE THR ASP PHE LEU ARG ARG GLN ALA PRO GLU LEU LEU PRO ALA ALA ILE SER SER GLY GLY ALA PRO LEU LEU GLY GLY ASP ALA GLN LEU LEU PRO HIS GLY T301 C482

S522  
GLY  
ALA  
ASP  
THR  
PHE  
GLY  
SER  
ASP  
GLY  
GLU  
LYS

● Molecule 3: Proteasome subunit beta



MET THR TRP PRO LEU LEU PRO ASP ARG LEU SER ILE ASN SER LEU SER GLY THR PRO PRO ALA VAL ASP LEU SER SER PHE THR ASP PHE LEU ARG ARG GLN ALA PRO GLU LEU LEU PRO ALA ALA ILE SER SER GLY GLY ALA PRO LEU LEU GLY GLY ASP ALA GLN LEU LEU PRO HIS GLY T301 S522

GLY  
ALA  
ASP  
THR  
PHE  
GLY  
SER  
ASP  
GLY  
GLU  
LYS

● Molecule 3: Proteasome subunit beta



MET THR TRP PRO LEU LEU PRO ASP ARG LEU SER ILE ASN SER LEU SER GLY THR PRO PRO ALA VAL ASP LEU SER SER PHE THR ASP PHE LEU ARG ARG GLN ALA PRO GLU LEU LEU PRO ALA ALA ILE SER SER GLY GLY ALA PRO LEU LEU GLY GLY ASP ALA GLN LEU LEU PRO HIS GLY T301 S522

GLY  
ALA  
ASP  
THR  
PHE  
GLY  
SER  
ASP  
GLY  
GLU  
LYS

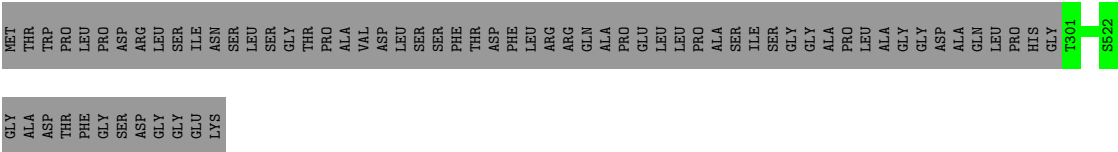
● Molecule 3: Proteasome subunit beta



MET THR TRP PRO LEU LEU PRO ASP ARG LEU SER ILE ASN SER LEU SER GLY THR PRO PRO ALA VAL ASP LEU SER SER PHE THR ASP PHE LEU ARG ARG GLN ALA PRO GLU LEU LEU PRO ALA ALA ILE SER SER GLY GLY ALA PRO LEU LEU GLY GLY ASP ALA GLN LEU LEU PRO HIS GLY T301 P309

S522  
GLY  
ALA  
ASP  
THR  
PHE  
GLY  
SER  
ASP  
GLY  
GLU  
LYS





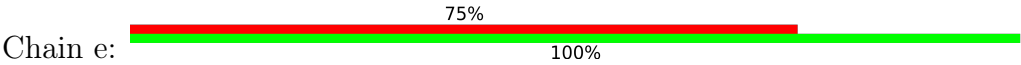
• Molecule 3: Proteasome subunit beta



• Molecule 4: Proteasome-associated ATPase



• Molecule 4: Proteasome-associated ATPase



• Molecule 4: Proteasome-associated ATPase



• Molecule 4: Proteasome-associated ATPase



• Molecule 4: Proteasome-associated ATPase





- Molecule 4: Proteasome-associated ATPase



- Molecule 4: Proteasome-associated ATPase





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	508000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	1.32	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	5000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	4.140	Depositor
Minimum map value	-1.162	Depositor
Average map value	-0.012	Depositor
Map value standard deviation	0.193	Depositor
Recommended contour level	0.6	Depositor
Map size ( $\text{\AA}$ )	423.936, 423.936, 423.936	wwPDB
Map dimensions	128, 128, 128	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	3.312, 3.312, 3.312	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	0.29	0/3808	0.52	0/5150
1	B	0.27	0/3766	0.53	0/5092
1	C	0.27	0/3706	0.51	0/5011
1	D	0.27	0/3651	0.52	0/4933
1	E	0.27	0/3655	0.52	0/4939
1	F	0.27	0/3646	0.52	0/4929
2	G	0.35	0/1683	0.56	0/2274
2	H	0.31	0/1683	0.57	0/2274
2	I	0.33	0/1683	0.56	0/2274
2	J	0.33	0/1683	0.55	0/2274
2	K	0.35	0/1683	0.56	0/2274
2	L	0.33	0/1683	0.56	0/2274
2	M	0.35	0/1683	0.56	0/2274
2	N	0.30	0/1683	0.54	0/2274
2	O	0.35	0/1683	0.55	0/2274
2	k	0.32	0/1683	0.55	0/2274
2	l	0.36	0/1683	0.57	0/2274
2	m	0.38	0/1683	0.56	0/2274
2	n	0.31	0/1683	0.55	0/2274
2	o	0.33	0/1683	0.56	0/2274
3	P	0.39	0/1662	0.53	0/2254
3	Q	0.30	0/1662	0.51	0/2254
3	R	0.32	0/1662	0.53	0/2254
3	S	0.32	0/1662	0.55	0/2254
3	T	0.30	0/1660	0.51	0/2251
3	U	0.30	0/1662	0.52	0/2254
3	V	0.33	0/1662	0.52	0/2254
3	W	0.35	0/1662	0.54	0/2254
3	X	0.30	0/1662	0.52	0/2254
3	Y	0.31	0/1662	0.54	0/2254
3	Z	0.37	0/1662	0.53	0/2254
3	a	0.34	0/1662	0.54	0/2254
3	b	0.32	0/1662	0.53	0/2254
3	c	0.33	0/1662	0.52	0/2254

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
4	d	0.24	0/34	0.35	0/43
4	e	0.28	0/34	0.32	0/43
4	f	0.24	0/34	0.49	0/43
4	g	0.27	0/34	0.55	0/43
4	h	0.26	0/34	0.53	0/43
4	i	0.26	0/34	0.47	0/43
4	j	0.28	0/34	0.50	0/43
All	All	0.31	0/69298	0.54	0/93744

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](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	474/609 (78%)	440 (93%)	33 (7%)	1 (0%)	47	81
1	B	467/609 (77%)	436 (93%)	30 (6%)	1 (0%)	47	81
1	C	462/609 (76%)	441 (96%)	18 (4%)	3 (1%)	25	66
1	D	450/609 (74%)	435 (97%)	15 (3%)	0	100	100
1	E	451/609 (74%)	436 (97%)	15 (3%)	0	100	100
1	F	452/609 (74%)	423 (94%)	28 (6%)	1 (0%)	47	81
2	G	211/248 (85%)	205 (97%)	6 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	211/248 (85%)	204 (97%)	7 (3%)	0	100	100
2	I	211/248 (85%)	202 (96%)	9 (4%)	0	100	100
2	J	211/248 (85%)	204 (97%)	7 (3%)	0	100	100
2	K	211/248 (85%)	206 (98%)	5 (2%)	0	100	100
2	L	211/248 (85%)	203 (96%)	8 (4%)	0	100	100
2	M	211/248 (85%)	202 (96%)	9 (4%)	0	100	100
2	N	211/248 (85%)	201 (95%)	10 (5%)	0	100	100
2	O	211/248 (85%)	203 (96%)	8 (4%)	0	100	100
2	k	211/248 (85%)	205 (97%)	6 (3%)	0	100	100
2	l	211/248 (85%)	204 (97%)	7 (3%)	0	100	100
2	m	211/248 (85%)	204 (97%)	7 (3%)	0	100	100
2	n	211/248 (85%)	204 (97%)	7 (3%)	0	100	100
2	o	211/248 (85%)	205 (97%)	6 (3%)	0	100	100
3	P	220/291 (76%)	212 (96%)	8 (4%)	0	100	100
3	Q	220/291 (76%)	215 (98%)	5 (2%)	0	100	100
3	R	220/291 (76%)	217 (99%)	3 (1%)	0	100	100
3	S	220/291 (76%)	211 (96%)	8 (4%)	1 (0%)	29	69
3	T	220/291 (76%)	215 (98%)	5 (2%)	0	100	100
3	U	220/291 (76%)	214 (97%)	6 (3%)	0	100	100
3	V	220/291 (76%)	217 (99%)	2 (1%)	1 (0%)	29	69
3	W	220/291 (76%)	211 (96%)	9 (4%)	0	100	100
3	X	220/291 (76%)	212 (96%)	8 (4%)	0	100	100
3	Y	220/291 (76%)	216 (98%)	4 (2%)	0	100	100
3	Z	220/291 (76%)	214 (97%)	6 (3%)	0	100	100
3	a	220/291 (76%)	215 (98%)	5 (2%)	0	100	100
3	b	220/291 (76%)	218 (99%)	2 (1%)	0	100	100
3	c	220/291 (76%)	214 (97%)	6 (3%)	0	100	100
4	d	2/4 (50%)	2 (100%)	0	0	100	100
4	e	2/4 (50%)	2 (100%)	0	0	100	100
4	f	2/4 (50%)	2 (100%)	0	0	100	100
4	g	2/4 (50%)	2 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	h	2/4 (50%)	2 (100%)	0	0	100	100
4	i	2/4 (50%)	2 (100%)	0	0	100	100
4	j	2/4 (50%)	2 (100%)	0	0	100	100
All	All	8804/11228 (78%)	8478 (96%)	318 (4%)	8 (0%)	54	86

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	146	VAL
1	C	222	ARG
3	S	482	GLY
1	C	217	LYS
1	B	150	VAL

### 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	405/511 (79%)	402 (99%)	3 (1%)	84	90
1	B	402/511 (79%)	399 (99%)	3 (1%)	84	90
1	C	389/511 (76%)	386 (99%)	3 (1%)	81	89
1	D	389/511 (76%)	388 (100%)	1 (0%)	92	95
1	E	389/511 (76%)	386 (99%)	3 (1%)	81	89
1	F	389/511 (76%)	387 (100%)	2 (0%)	88	93
2	G	165/192 (86%)	163 (99%)	2 (1%)	71	83
2	H	165/192 (86%)	163 (99%)	2 (1%)	71	83
2	I	165/192 (86%)	163 (99%)	2 (1%)	71	83
2	J	165/192 (86%)	164 (99%)	1 (1%)	86	92
2	K	165/192 (86%)	165 (100%)	0	100	100
2	L	165/192 (86%)	163 (99%)	2 (1%)	71	83
2	M	165/192 (86%)	162 (98%)	3 (2%)	59	77

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	N	165/192 (86%)	164 (99%)	1 (1%)	86	92
2	O	165/192 (86%)	165 (100%)	0	100	100
2	k	165/192 (86%)	164 (99%)	1 (1%)	86	92
2	l	165/192 (86%)	164 (99%)	1 (1%)	86	92
2	m	165/192 (86%)	165 (100%)	0	100	100
2	n	165/192 (86%)	164 (99%)	1 (1%)	86	92
2	o	165/192 (86%)	164 (99%)	1 (1%)	86	92
3	P	165/217 (76%)	165 (100%)	0	100	100
3	Q	165/217 (76%)	165 (100%)	0	100	100
3	R	165/217 (76%)	165 (100%)	0	100	100
3	S	165/217 (76%)	165 (100%)	0	100	100
3	T	164/217 (76%)	164 (100%)	0	100	100
3	U	165/217 (76%)	165 (100%)	0	100	100
3	V	165/217 (76%)	165 (100%)	0	100	100
3	W	165/217 (76%)	165 (100%)	0	100	100
3	X	165/217 (76%)	165 (100%)	0	100	100
3	Y	165/217 (76%)	165 (100%)	0	100	100
3	Z	165/217 (76%)	165 (100%)	0	100	100
3	a	165/217 (76%)	165 (100%)	0	100	100
3	b	165/217 (76%)	165 (100%)	0	100	100
3	c	165/217 (76%)	164 (99%)	1 (1%)	86	92
4	d	3/3 (100%)	3 (100%)	0	100	100
4	e	3/3 (100%)	3 (100%)	0	100	100
4	f	3/3 (100%)	3 (100%)	0	100	100
4	g	3/3 (100%)	3 (100%)	0	100	100
4	h	3/3 (100%)	3 (100%)	0	100	100
4	i	3/3 (100%)	3 (100%)	0	100	100
4	j	3/3 (100%)	3 (100%)	0	100	100
All	All	7003/8813 (80%)	6970 (100%)	33 (0%)	89	93

5 of 33 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	M	134	LYS
2	M	141	ILE
3	c	381	ASN
1	E	324	MET
1	E	254	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
2	H	98	GLN
2	L	165	ASN
3	a	437	GLN
2	M	114	GLN
3	S	396	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

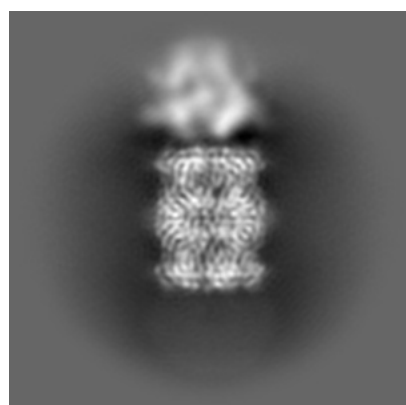
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-27226. These allow visual inspection of the internal detail of the map and identification of artifacts.

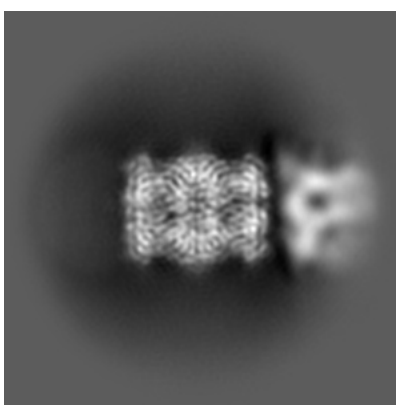
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

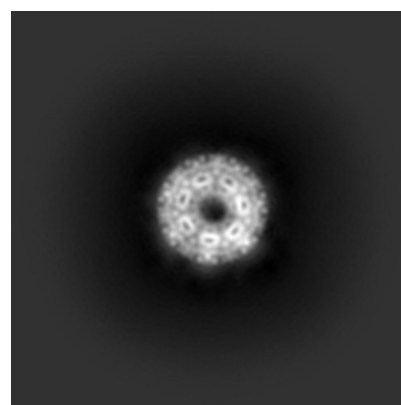
#### 6.1.1 Primary map



X



Y

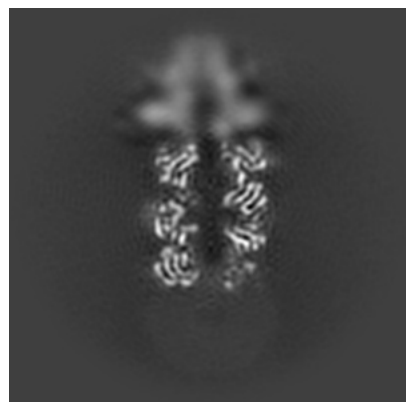


Z

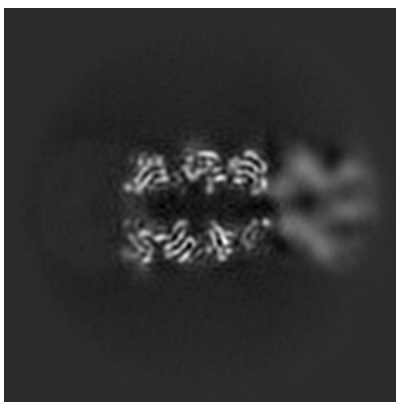
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

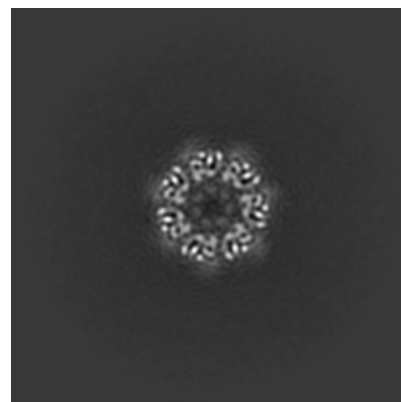
#### 6.2.1 Primary map



X Index: 64



Y Index: 64



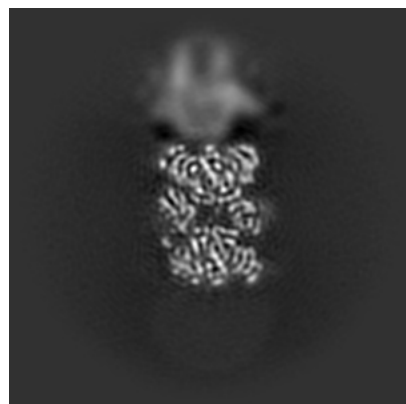
Z Index: 64



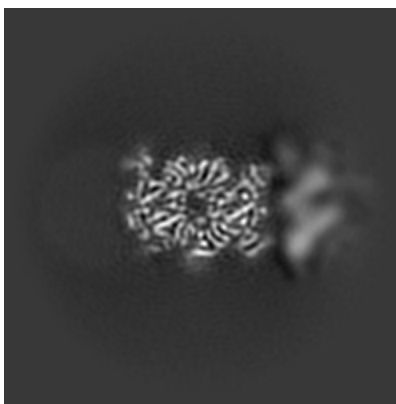
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

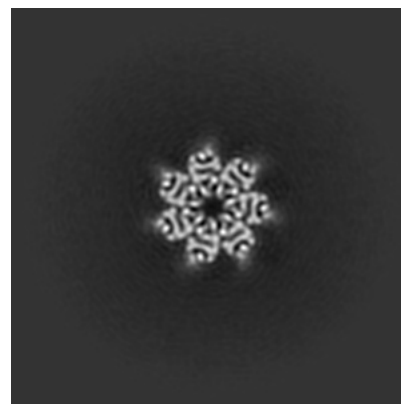
### 6.3.1 Primary map



X Index: 72



Y Index: 72



Z Index: 79

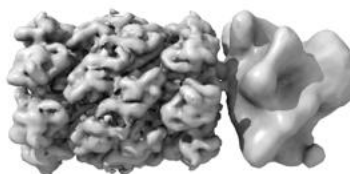
The images above show the largest variance slices of the map in three orthogonal directions.

## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.6. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

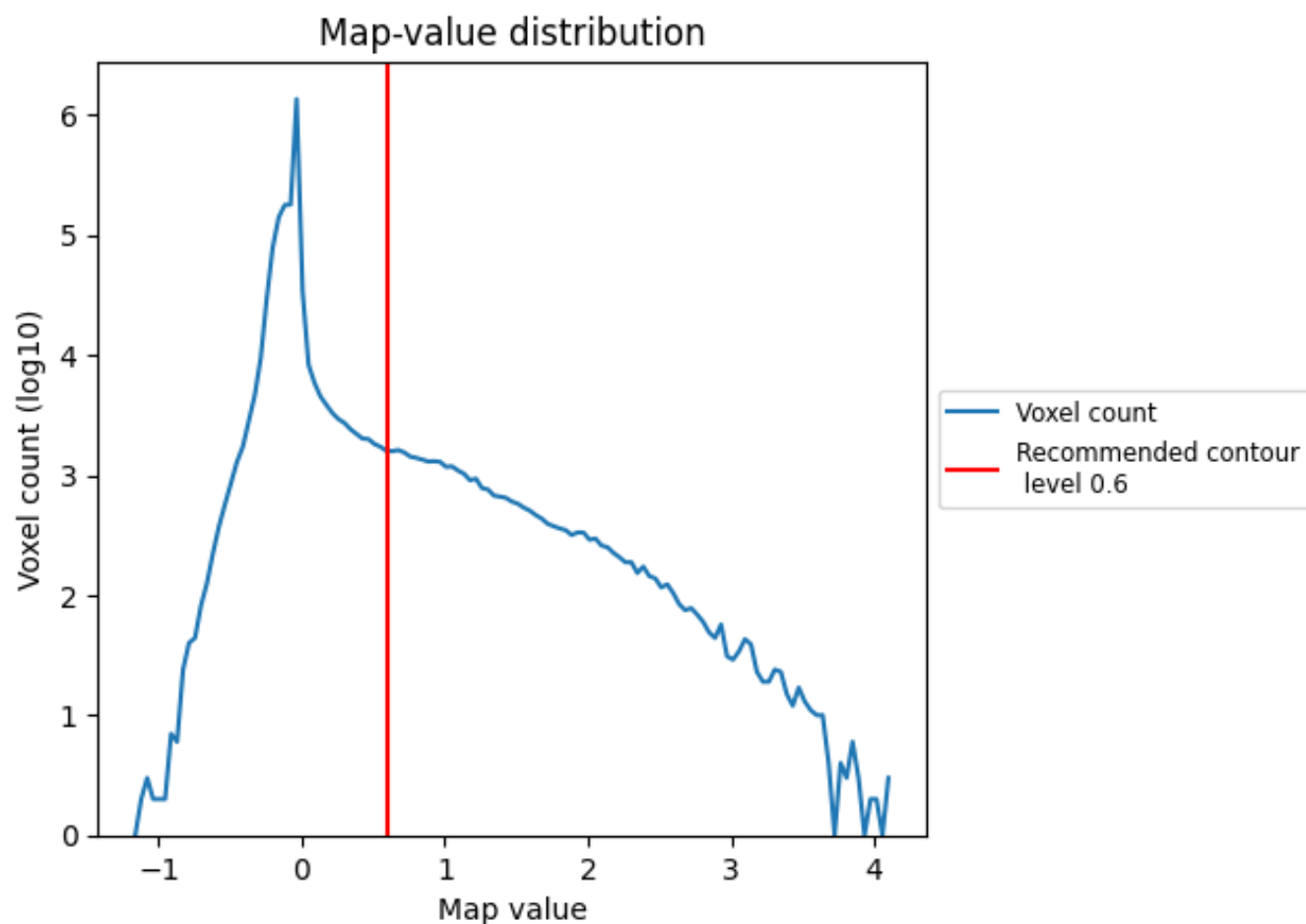
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

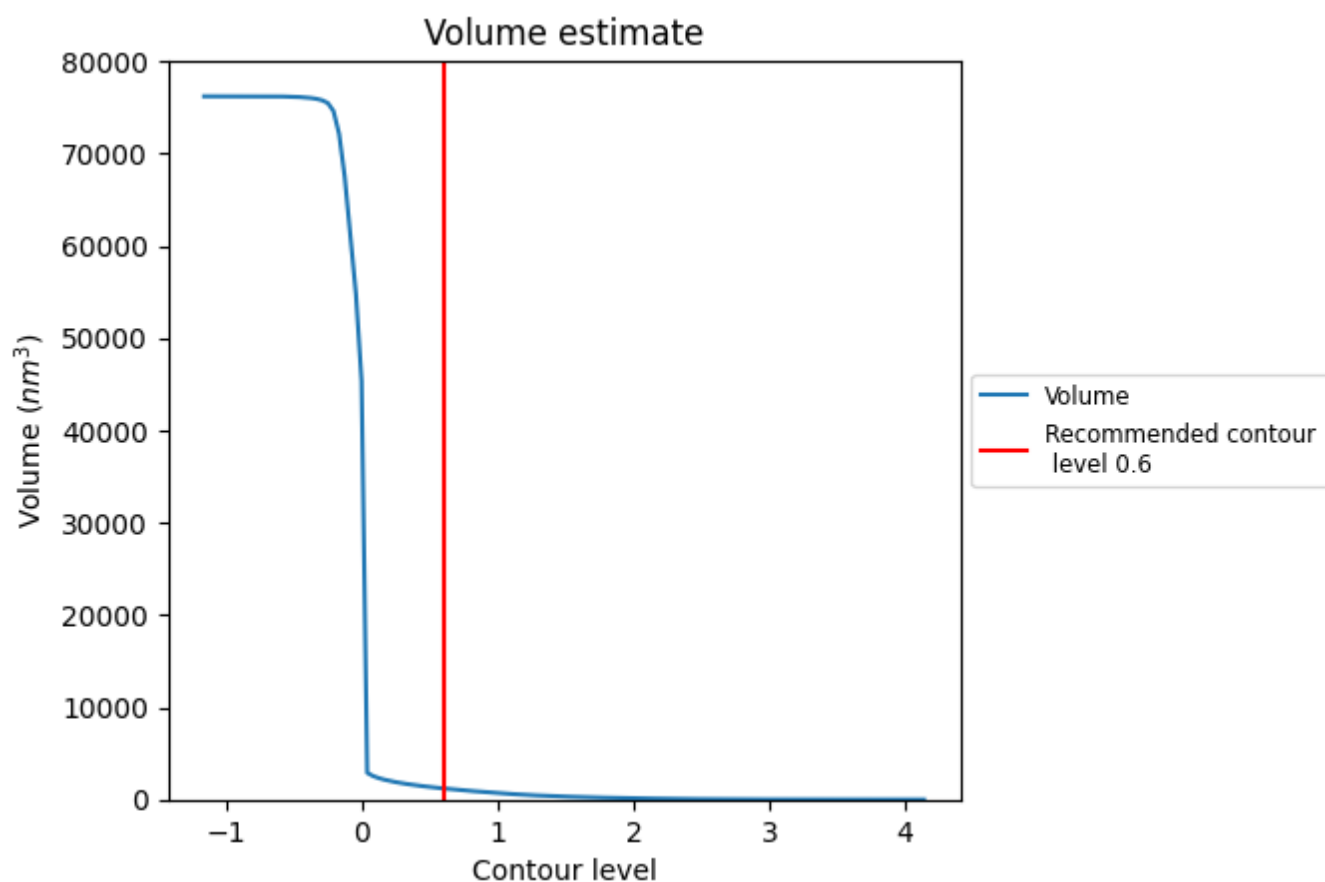
This section contains the results of statistical analysis of the map.

### 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.

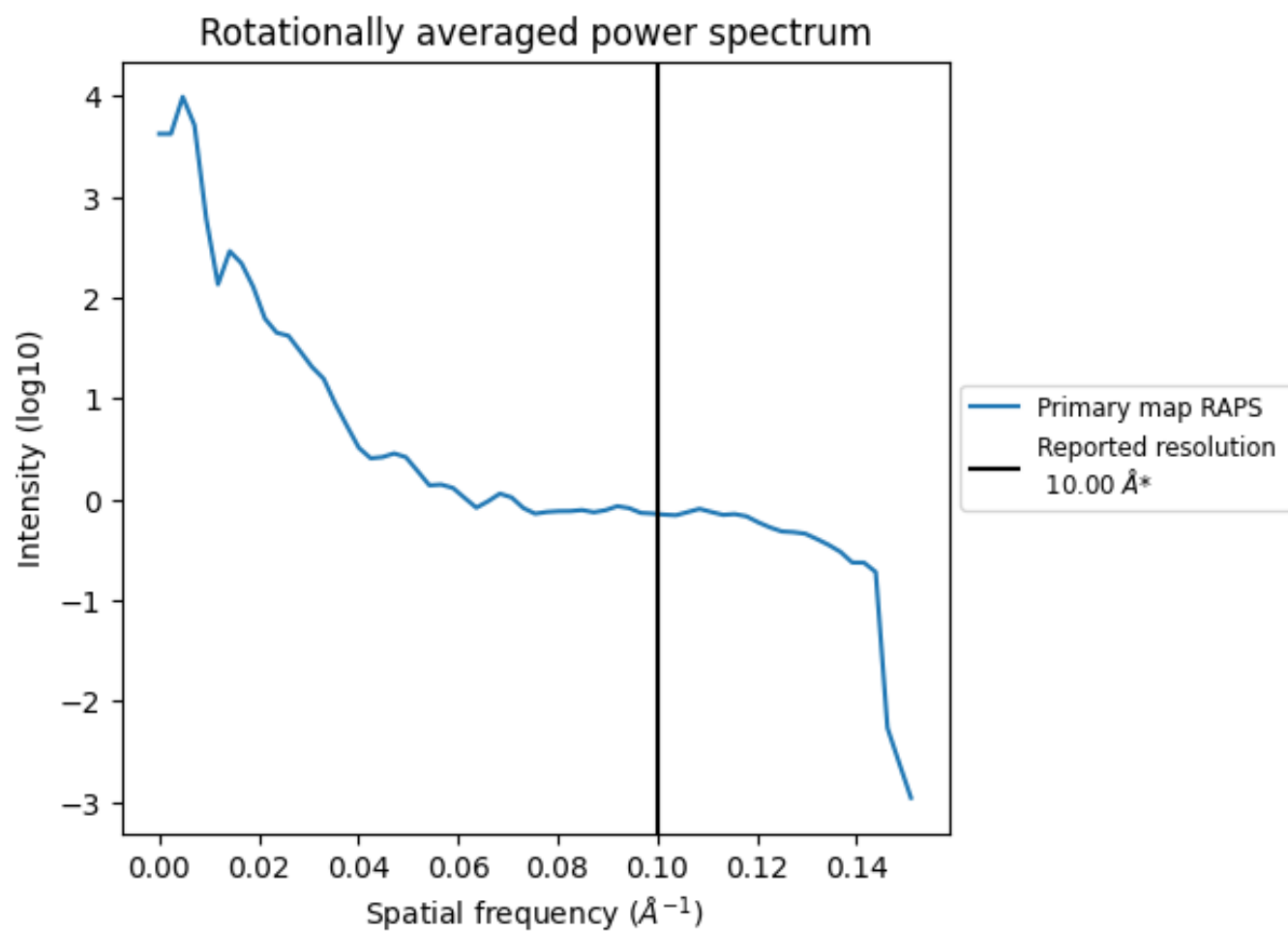
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1208 nm<sup>3</sup>; this corresponds to an approximate mass of 1092 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 0.100 Å<sup>-1</sup>

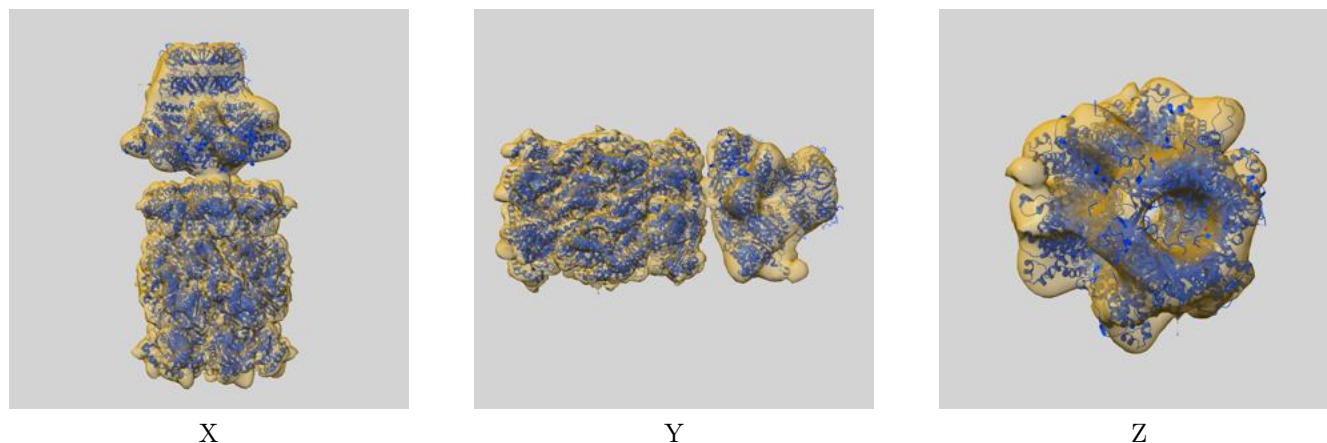
## 8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

## 9 Map-model fit [i](#)

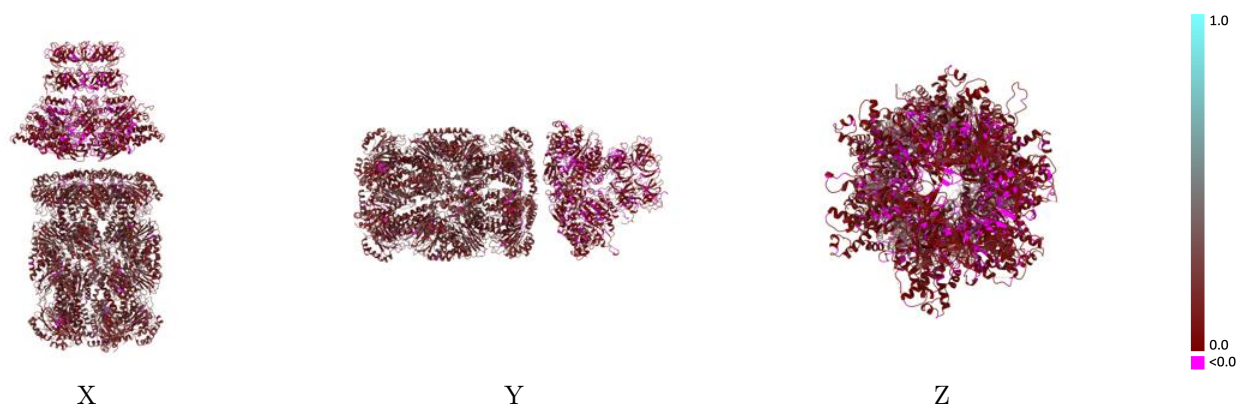
This section contains information regarding the fit between EMDB map EMD-27226 and PDB model 8D6Y. Per-residue inclusion information can be found in section [3](#) on page [8](#).

### 9.1 Map-model overlay [i](#)



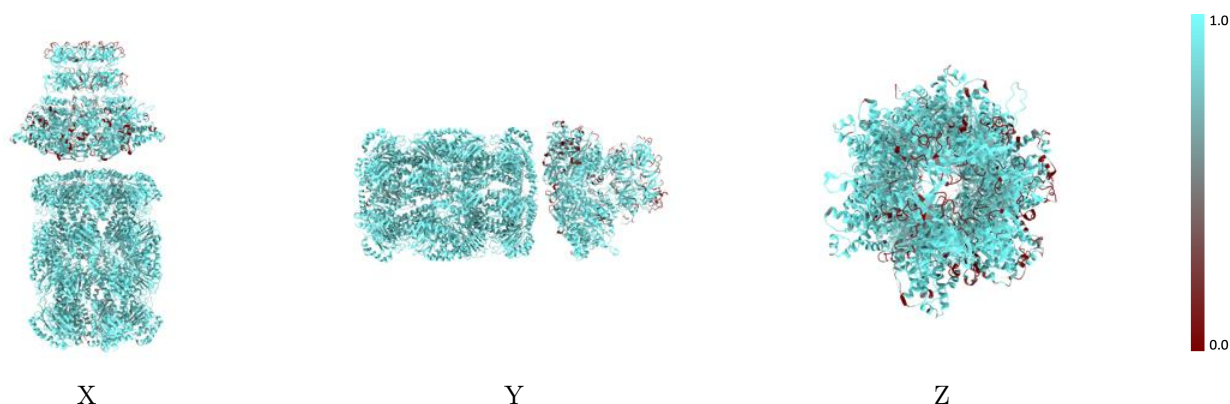
The images above show the 3D surface view of the map at the recommended contour level 0.6 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

## 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.

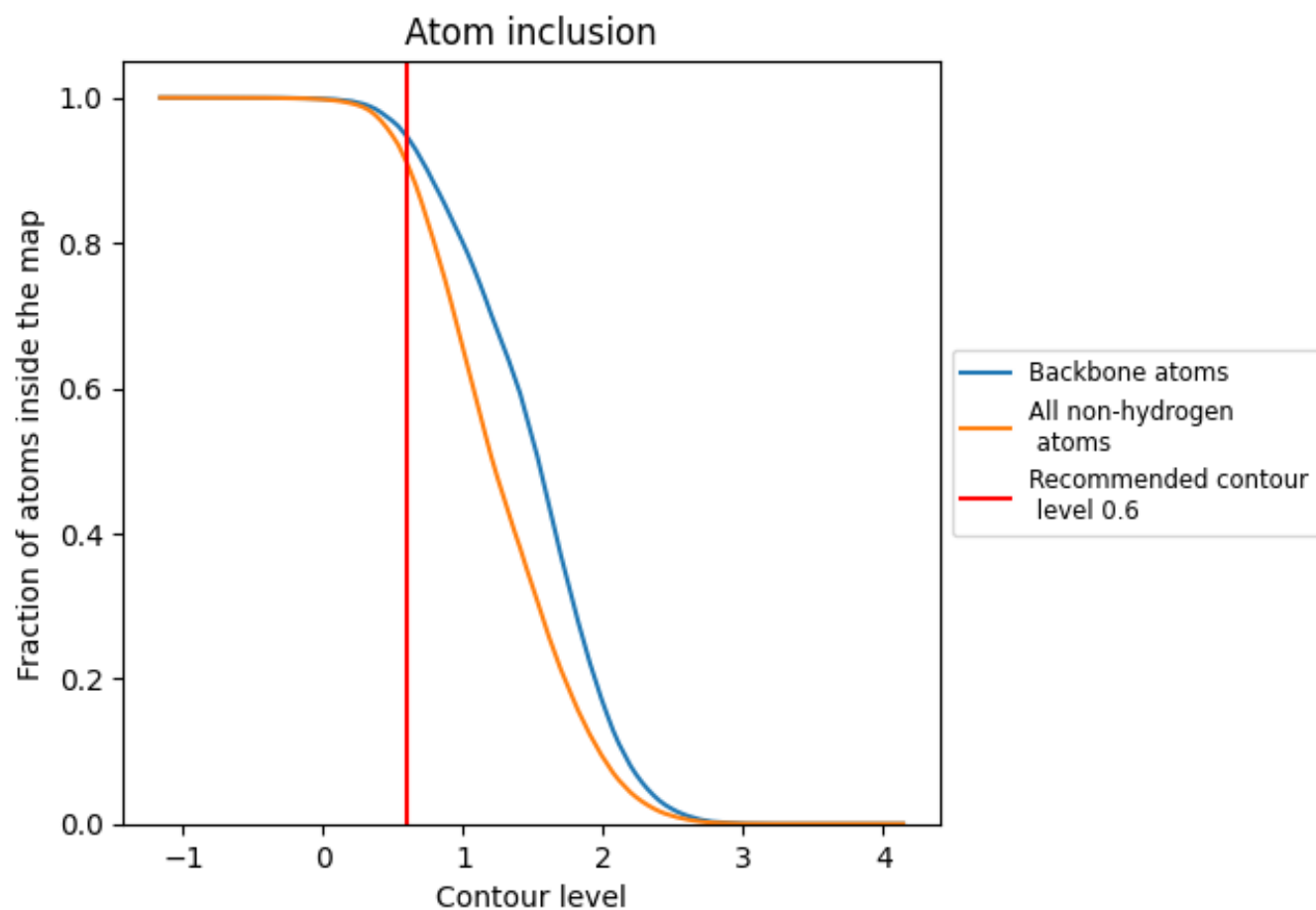
## 9.3 Atom inclusion mapped to coordinate model [i](#)



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 (0.6).









































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 91% 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 (0.6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9092	 0.1510
A	 0.8198	 0.0580
B	 0.8405	 0.0600
C	 0.8685	 0.0600
D	 0.8450	 0.0620
E	 0.8051	 0.0530
F	 0.7736	 0.0500
G	 0.9608	 0.1930
H	 0.9645	 0.1900
I	 0.9626	 0.1950
J	 0.9328	 0.1830
K	 0.9396	 0.1900
L	 0.9228	 0.1900
M	 0.9247	 0.1860
N	 0.9340	 0.1910
O	 0.9253	 0.1870
P	 0.9625	 0.2000
Q	 0.9625	 0.2090
R	 0.9644	 0.2050
S	 0.9594	 0.2070
T	 0.9550	 0.2060
U	 0.9650	 0.2100
V	 0.9588	 0.2010
W	 0.9438	 0.1990
X	 0.9613	 0.2100
Y	 0.9513	 0.1980
Z	 0.9569	 0.2040
a	 0.9532	 0.1980
b	 0.9507	 0.1950
c	 0.9501	 0.2050
d	 0.5152	 -0.1190
e	 0.3030	 -0.1590
f	 0.3333	 -0.1640
g	 0.5455	 -0.1860
h	 0.3939	 -0.1990



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Chain	Atom inclusion	Q-score
i	 0.5152	 -0.0840
j	 0.6970	 -0.0030
k	 0.9328	 0.1950
l	 0.9601	 0.1930
m	 0.9533	 0.1920
n	 0.9658	 0.1910
o	 0.9626	 0.1920