



wwPDB EM Validation Summary Report ⓘ

Jul 7, 2024 – 10:22 pm BST

PDB ID : 7PG9
EMDB ID : EMD-13389
Title : human 20S proteasome
Authors : Xu, C.; Cong, Y.
Deposited on : 2021-08-13
Resolution : 3.70 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.dev92
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

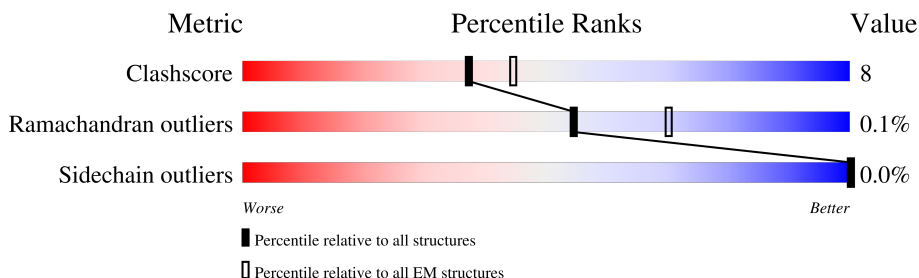
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.70 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.






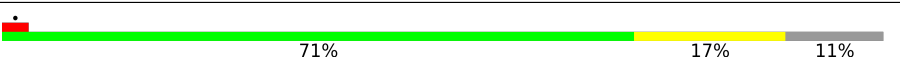
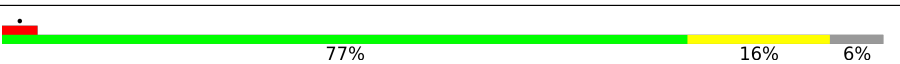

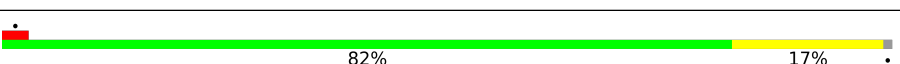
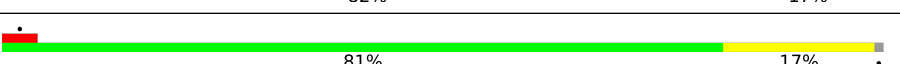
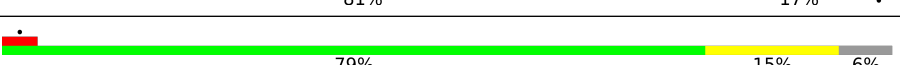
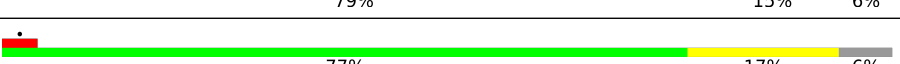
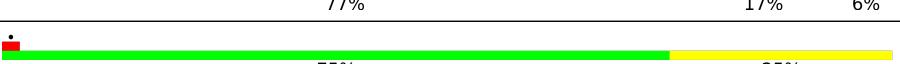
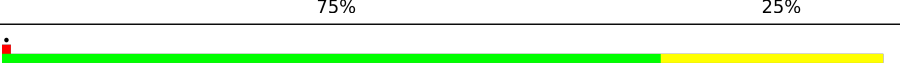

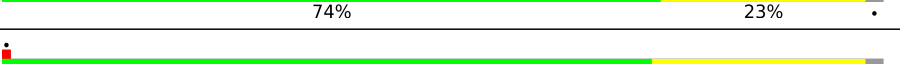



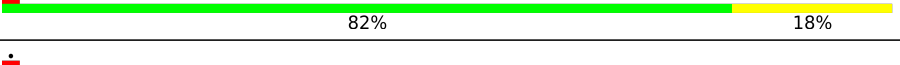
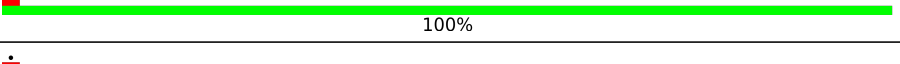

Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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 ≥ 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	246	
1	O	246	
2	B	234	
2	P	234	
3	C	261	
3	Q	261	
4	D	248	
4	R	248	

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Mol	Chain	Length	Quality of chain
5	E	241	
5	S	241	
6	F	263	
6	T	263	
7	G	255	
7	U	255	
8	H	205	
8	V	205	
9	I	234	
9	W	234	
10	J	205	
10	X	205	
11	K	201	
11	Y	201	
12	L	204	
12	Z	204	
13	M	213	
13	a	213	
14	N	219	
14	b	219	

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 45326 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 type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	239	Total	C	N	O	S	0	0
			1729	1101	303	313	12		
1	O	238	Total	C	N	O	S	0	0
			1725	1099	302	312	12		

- Molecule 2 is a protein called Proteasome subunit alpha type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	229	Total	C	N	O	S	0	0
			1662	1080	288	288	6		
2	P	229	Total	C	N	O	S	0	0
			1662	1080	288	288	6		

- Molecule 3 is a protein called Proteasome subunit alpha type-4.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	247	Total	C	N	O	S	0	0
			1786	1143	320	313	10		
3	Q	247	Total	C	N	O	S	0	0
			1786	1143	320	313	10		

- Molecule 4 is a protein called Proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	232	Total	C	N	O	S	0	0
			1633	1038	306	284	5		
4	R	232	Total	C	N	O	S	0	0
			1633	1038	306	284	5		

- Molecule 5 is a protein called Proteasome subunit alpha type-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	233	Total	C	N	O	S	0	0
			1659	1056	287	305	11		
5	S	233	Total	C	N	O	S	0	0
			1659	1056	287	305	11		

- Molecule 6 is a protein called Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	233	Total	C	N	O	S	0	0
			1704	1087	315	293	9		
6	T	233	Total	C	N	O	S	0	0
			1704	1087	315	293	9		

- Molecule 7 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	239	Total	C	N	O	S	0	0
			1760	1131	308	311	10		
7	U	239	Total	C	N	O	S	0	0
			1760	1131	308	311	10		

- Molecule 8 is a protein called Proteasome subunit beta type-6.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	202	Total	C	N	O	S	0	0
			1465	926	256	271	12		
8	V	202	Total	C	N	O	S	0	0
			1465	926	256	271	12		

- Molecule 9 is a protein called Proteasome subunit beta type-7.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	220	Total	C	N	O	S	0	0
			1576	1003	272	292	9		
9	W	220	Total	C	N	O	S	0	0
			1576	1003	272	292	9		

- Molecule 10 is a protein called Proteasome subunit beta type-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	204	Total	C	N	O	S	0	0
			1534	987	262	267	18		

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Mol	Chain	Residues	Atoms					AltConf	Trace
10	X	204	Total	C	N	O	S	0	0
			1534	987	262	267	18		

- Molecule 11 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	196	Total	C	N	O	S	0	0
			1506	973	259	266	8		
11	Y	196	Total	C	N	O	S	0	0
			1506	973	259	266	8		

- Molecule 12 is a protein called Proteasome subunit beta type-5.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	200	Total	C	N	O	S	0	0
			1500	954	270	267	9		
12	Z	200	Total	C	N	O	S	0	0
			1500	954	270	267	9		

- Molecule 13 is a protein called Proteasome subunit beta type-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	212	Total	C	N	O	S	0	0
			1583	1016	279	278	10		
13	a	212	Total	C	N	O	S	0	0
			1583	1016	279	278	10		

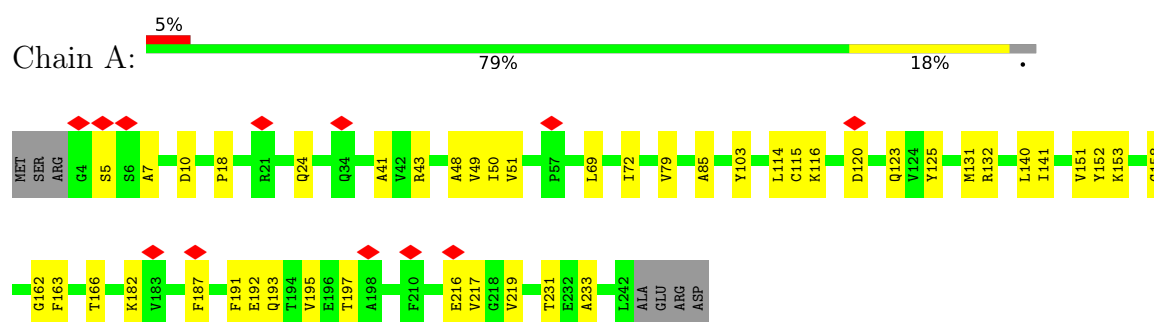
- Molecule 14 is a protein called Proteasome subunit beta type-4.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	212	Total	C	N	O	S	0	0
			1568	998	279	280	11		
14	b	212	Total	C	N	O	S	0	0
			1568	998	279	280	11		

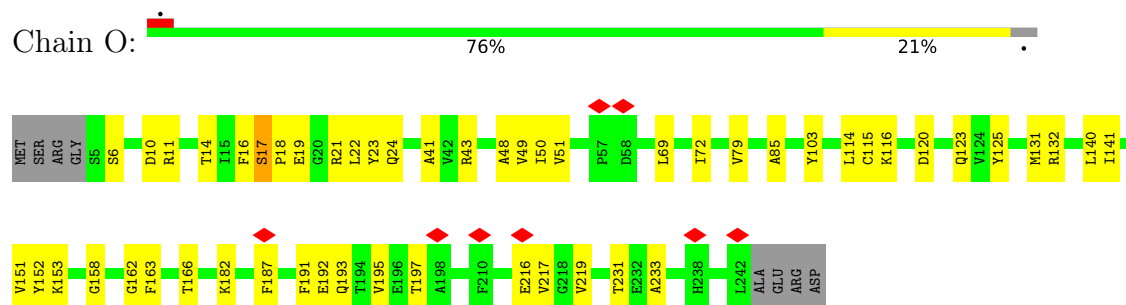
3 Residue-property plots [i](#)

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

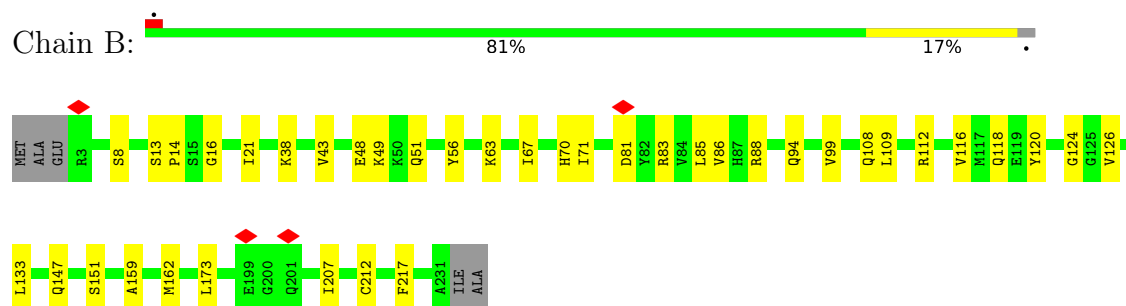
- Molecule 1: Proteasome subunit alpha type-6



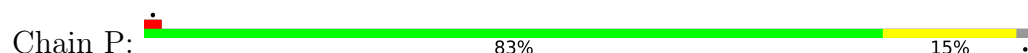
- Molecule 1: Proteasome subunit alpha type-6



- Molecule 2: Proteasome subunit alpha type-2



- Molecule 2: Proteasome subunit alpha type-2

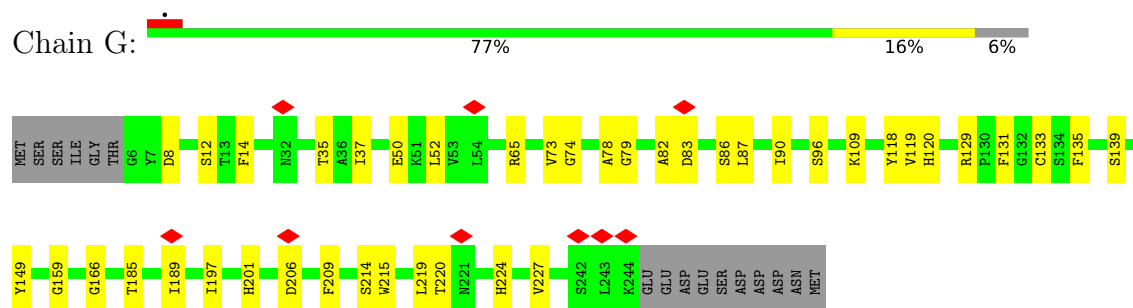




HIS

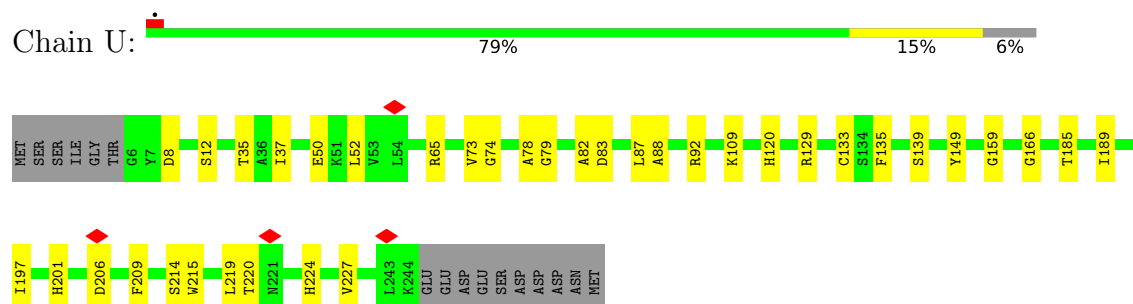
- Molecule 7: Proteasome subunit alpha type-3

Chain G:



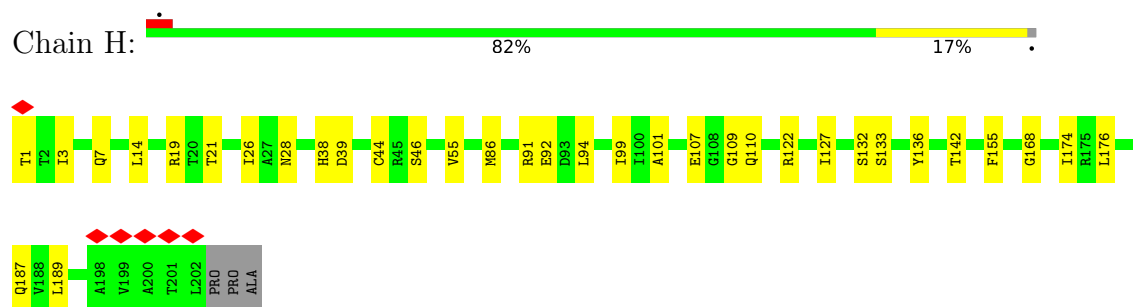
- Molecule 7: Proteasome subunit alpha type-3

Chain U:



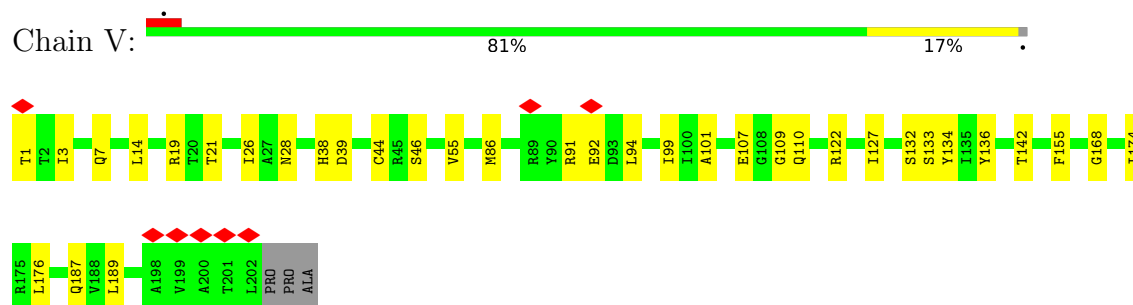
- Molecule 8: Proteasome subunit beta type-6

Chain H:



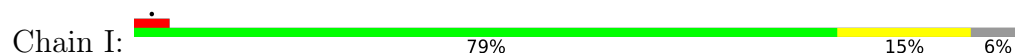
- Molecule 8: Proteasome subunit beta type-6

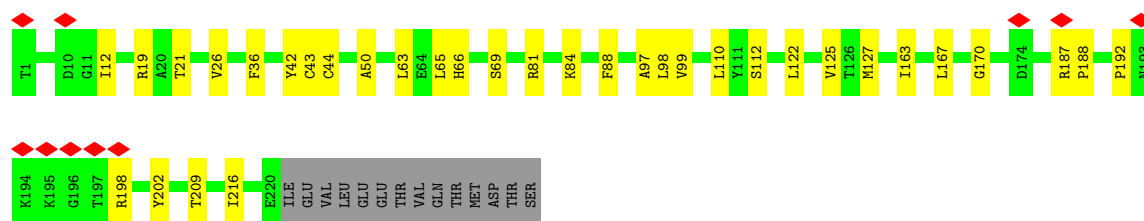
Chain V:



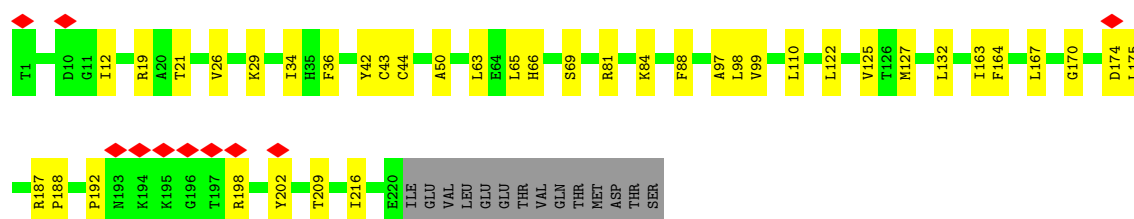
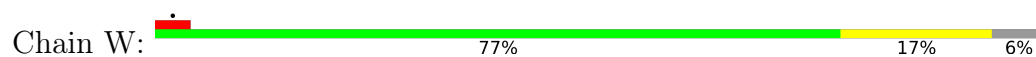
- Molecule 9: Proteasome subunit beta type-7

Chain I:

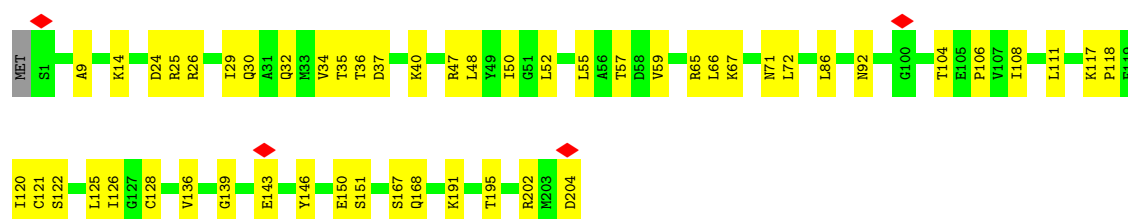




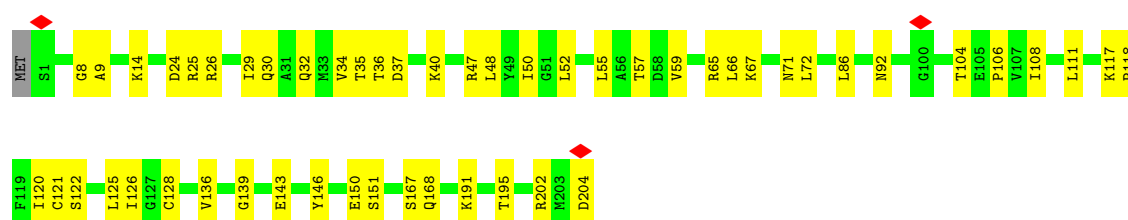
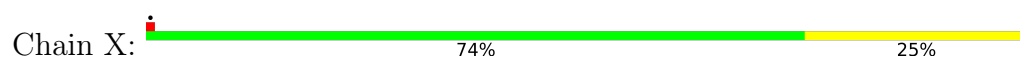
- Molecule 9: Proteasome subunit beta type-7



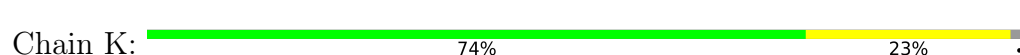
- Molecule 10: Proteasome subunit beta type-3



- Molecule 10: Proteasome subunit beta type-3



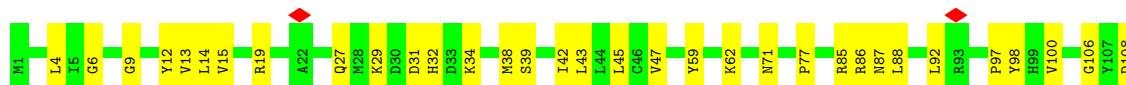
- Molecule 11: Proteasome subunit beta type-2





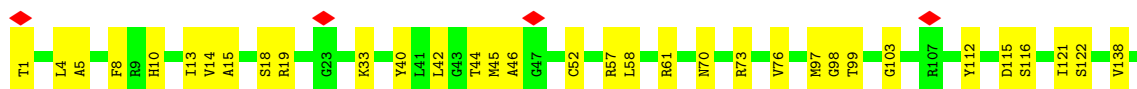
- Molecule 11: Proteasome subunit beta type-2

Chain Y: 73% 24%



- Molecule 12: Proteasome subunit beta type-5

Chain L: 77% 21%



- Molecule 12: Proteasome subunit beta type-5

Chain Z: 77% 21%

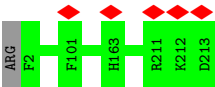


- Molecule 13: Proteasome subunit beta type-1

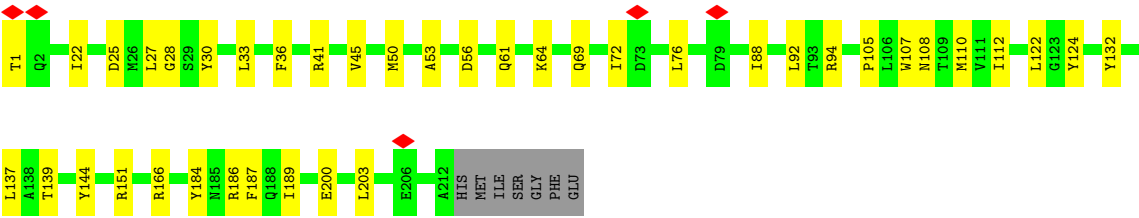
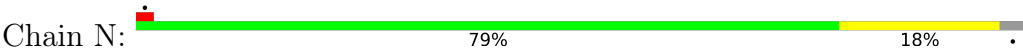
Chain M: 82% 18%



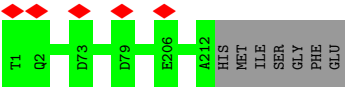
- Molecule 13: Proteasome subunit beta type-1



• Molecule 14: Proteasome subunit beta type-4



• Molecule 14: Proteasome subunit beta type-4



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	154436	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 ($e^-/\text{\AA}^2$)	38	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.058	Depositor
Minimum map value	-0.022	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0153	Depositor
Map size (\AA)	289.96, 289.96, 289.96	wwPDB
Map dimensions	220, 220, 220	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.318, 1.318, 1.318	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.26	0/1758	0.44	0/2386
1	O	0.27	0/1754	0.44	0/2381
2	B	0.26	0/1701	0.42	0/2318
2	P	0.25	0/1701	0.42	0/2318
3	C	0.24	0/1815	0.43	0/2466
3	Q	0.24	0/1815	0.43	0/2466
4	D	0.24	0/1657	0.44	0/2261
4	R	0.24	0/1657	0.43	0/2261
5	E	0.25	0/1685	0.43	0/2290
5	S	0.24	0/1685	0.43	0/2290
6	F	0.24	0/1738	0.43	0/2364
6	T	0.24	0/1738	0.43	0/2364
7	G	0.26	0/1795	0.42	0/2434
7	U	0.26	0/1795	0.42	0/2434
8	H	0.25	0/1491	0.43	0/2021
8	V	0.25	0/1491	0.43	0/2021
9	I	0.25	0/1603	0.45	0/2180
9	W	0.25	0/1603	0.45	0/2180
10	J	0.25	0/1563	0.45	0/2113
10	X	0.25	0/1563	0.45	0/2113
11	K	0.25	0/1538	0.46	0/2088
11	Y	0.25	0/1538	0.45	0/2088
12	L	0.25	0/1531	0.44	0/2076
12	Z	0.25	0/1531	0.44	0/2076
13	M	0.25	0/1613	0.45	0/2178
13	a	0.25	0/1613	0.45	0/2178
14	N	0.25	0/1598	0.45	0/2170
14	b	0.25	0/1598	0.45	0/2170
All	All	0.25	0/46168	0.44	0/62685

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

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	1729	0	1650	28	0
1	O	1725	0	1647	40	0
2	B	1662	0	1590	28	0
2	P	1662	0	1590	25	0
3	C	1786	0	1717	28	0
3	Q	1786	0	1717	27	0
4	D	1633	0	1518	31	0
4	R	1633	0	1518	29	0
5	E	1659	0	1589	30	0
5	S	1659	0	1589	32	0
6	F	1704	0	1638	39	0
6	T	1704	0	1638	35	0
7	G	1760	0	1680	31	0
7	U	1760	0	1680	27	0
8	H	1465	0	1419	20	0
8	V	1465	0	1419	22	0
9	I	1576	0	1558	29	0
9	W	1576	0	1558	35	0
10	J	1534	0	1539	34	0
10	X	1534	0	1539	36	0
11	K	1506	0	1475	41	0
11	Y	1506	0	1475	40	0
12	L	1500	0	1441	30	0
12	Z	1500	0	1441	32	0
13	M	1583	0	1579	29	0
13	a	1583	0	1579	0	0
14	N	1568	0	1513	28	0
14	b	1568	0	1513	0	0
All	All	45326	0	43809	679	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 8.

The worst 5 of 679 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)
9:W:187:ARG:HB3	9:W:188:PRO:HD3	1.24	1.19
9:I:187:ARG:HB2	9:I:188:PRO:HD3	1.26	1.16
5:S:121:LEU:HD12	6:T:79:ALA:HB3	1.59	0.84
5:E:121:LEU:HD12	6:F:79:ALA:HB3	1.59	0.83
9:I:187:ARG:CB	9:I:188:PRO:HD3	2.09	0.83

There are no symmetry-related clashes.

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	237/246 (96%)	229 (97%)	7 (3%)	1 (0%)	34	69
1	O	236/246 (96%)	229 (97%)	6 (2%)	1 (0%)	34	69
2	B	227/234 (97%)	222 (98%)	5 (2%)	0	100	100
2	P	227/234 (97%)	222 (98%)	5 (2%)	0	100	100
3	C	245/261 (94%)	231 (94%)	14 (6%)	0	100	100
3	Q	245/261 (94%)	231 (94%)	14 (6%)	0	100	100
4	D	230/248 (93%)	221 (96%)	8 (4%)	1 (0%)	34	69
4	R	230/248 (93%)	223 (97%)	6 (3%)	1 (0%)	34	69
5	E	231/241 (96%)	226 (98%)	5 (2%)	0	100	100
5	S	231/241 (96%)	226 (98%)	5 (2%)	0	100	100
6	F	231/263 (88%)	227 (98%)	3 (1%)	1 (0%)	34	69
6	T	231/263 (88%)	227 (98%)	3 (1%)	1 (0%)	34	69
7	G	237/255 (93%)	232 (98%)	5 (2%)	0	100	100
7	U	237/255 (93%)	232 (98%)	5 (2%)	0	100	100
8	H	200/205 (98%)	197 (98%)	3 (2%)	0	100	100
8	V	200/205 (98%)	197 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	I	218/234 (93%)	204 (94%)	14 (6%)	0	100	100
9	W	218/234 (93%)	204 (94%)	14 (6%)	0	100	100
10	J	202/205 (98%)	194 (96%)	8 (4%)	0	100	100
10	X	202/205 (98%)	194 (96%)	8 (4%)	0	100	100
11	K	194/201 (96%)	189 (97%)	5 (3%)	0	100	100
11	Y	194/201 (96%)	189 (97%)	5 (3%)	0	100	100
12	L	198/204 (97%)	194 (98%)	4 (2%)	0	100	100
12	Z	198/204 (97%)	194 (98%)	4 (2%)	0	100	100
13	M	210/213 (99%)	204 (97%)	6 (3%)	0	100	100
13	a	210/213 (99%)	204 (97%)	6 (3%)	0	100	100
14	N	210/219 (96%)	205 (98%)	5 (2%)	0	100	100
14	b	210/219 (96%)	205 (98%)	5 (2%)	0	100	100
All	All	6139/6458 (95%)	5952 (97%)	181 (3%)	6 (0%)	54	83

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	R	216	SER
1	A	187	PHE
1	O	187	PHE
4	D	215	GLN
6	F	59	HIS

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	163/210 (78%)	163 (100%)	0	100	100
1	O	163/210 (78%)	161 (99%)	2 (1%)	71	84
2	B	150/191 (78%)	150 (100%)	0	100	100
2	P	150/191 (78%)	150 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	160/221 (72%)	160 (100%)	0	100	100
3	Q	160/221 (72%)	160 (100%)	0	100	100
4	D	136/211 (64%)	136 (100%)	0	100	100
4	R	136/211 (64%)	136 (100%)	0	100	100
5	E	158/203 (78%)	158 (100%)	0	100	100
5	S	158/203 (78%)	158 (100%)	0	100	100
6	F	160/224 (71%)	160 (100%)	0	100	100
6	T	160/224 (71%)	160 (100%)	0	100	100
7	G	162/212 (76%)	162 (100%)	0	100	100
7	U	162/212 (76%)	162 (100%)	0	100	100
8	H	140/159 (88%)	140 (100%)	0	100	100
8	V	140/159 (88%)	140 (100%)	0	100	100
9	I	157/195 (80%)	157 (100%)	0	100	100
9	W	157/195 (80%)	157 (100%)	0	100	100
10	J	155/174 (89%)	155 (100%)	0	100	100
10	X	155/174 (89%)	155 (100%)	0	100	100
11	K	148/171 (86%)	148 (100%)	0	100	100
11	Y	148/171 (86%)	148 (100%)	0	100	100
12	L	138/159 (87%)	138 (100%)	0	100	100
12	Z	138/159 (87%)	138 (100%)	0	100	100
13	M	158/178 (89%)	158 (100%)	0	100	100
13	a	158/178 (89%)	158 (100%)	0	100	100
14	N	149/181 (82%)	149 (100%)	0	100	100
14	b	149/181 (82%)	149 (100%)	0	100	100
All	All	4268/5378 (79%)	4266 (100%)	2 (0%)	100	100

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	O	17	SER
1	O	23	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 58 such sidechains are listed below:

Mol	Chain	Res	Type
1	O	128	ASN
13	a	157	ASN
3	Q	167	ASN
13	a	108	ASN
11	Y	32	HIS

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.

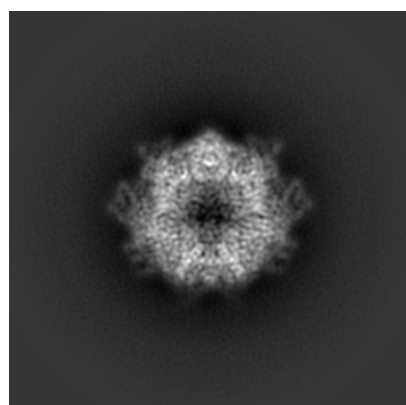
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13389. These allow visual inspection of the internal detail of the map and identification of artifacts.

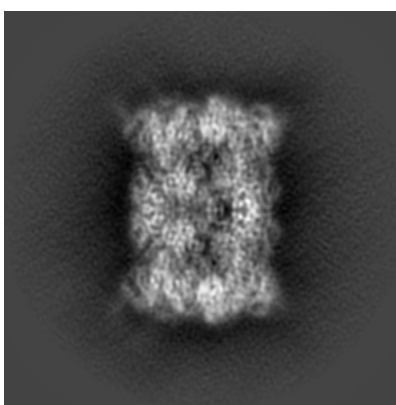
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

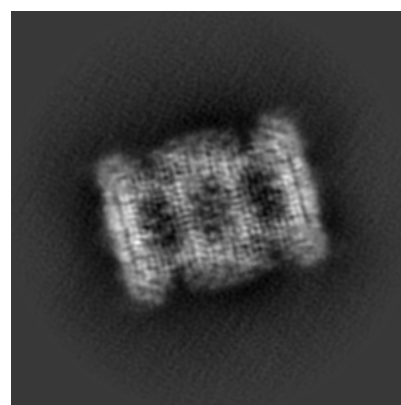
6.1.1 Primary map



X



Y

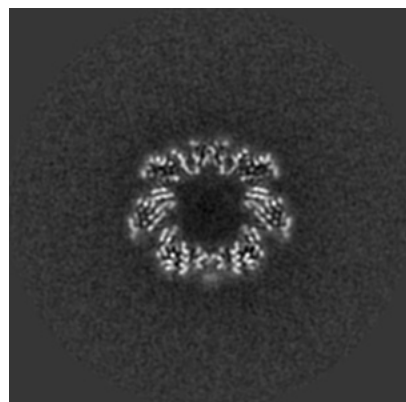


Z

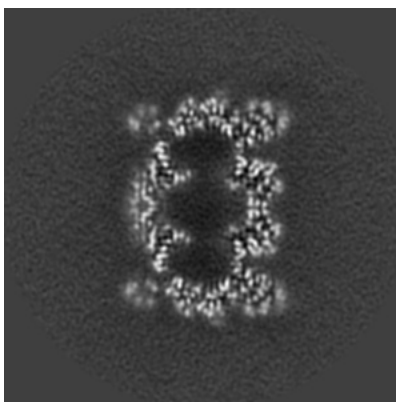
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

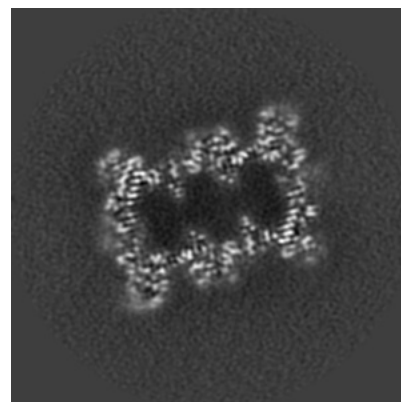
6.2.1 Primary map



X Index: 110



Y Index: 110

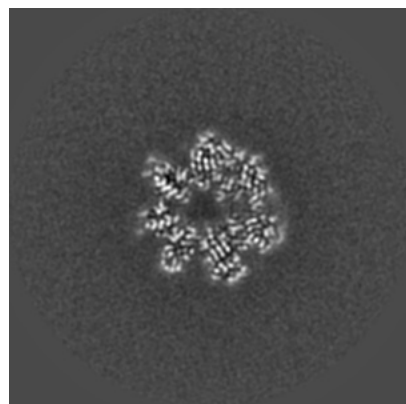


Z Index: 110

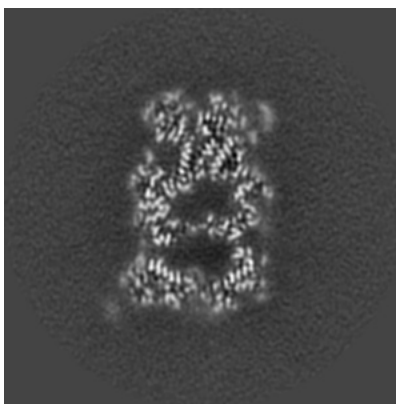
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

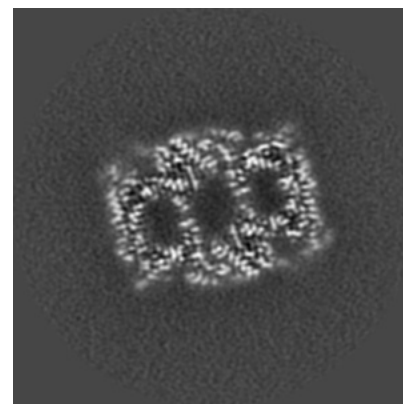
6.3.1 Primary map



X Index: 95



Y Index: 96

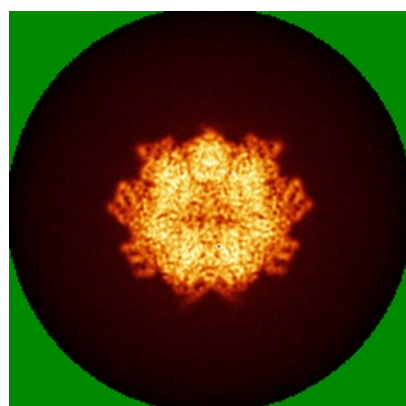


Z Index: 97

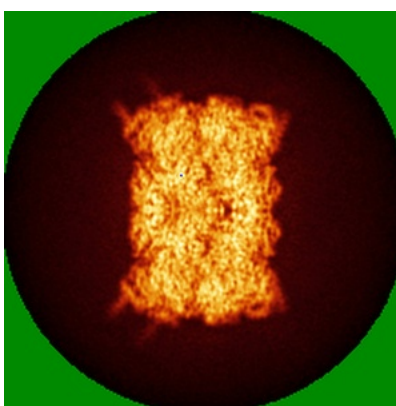
The images above show the largest variance slices of the map in three orthogonal directions.

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

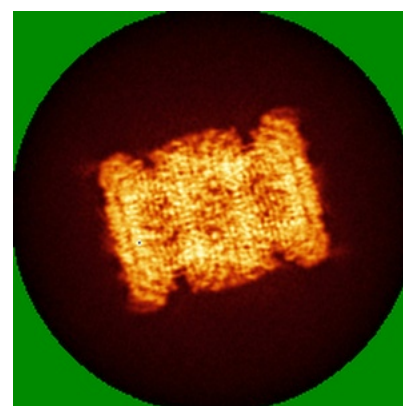
6.4.1 Primary map



X



Y

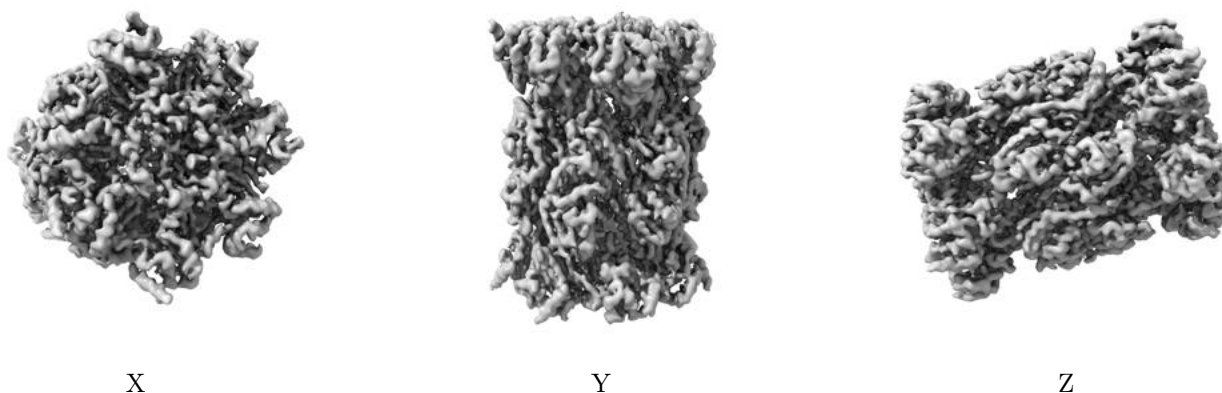


Z

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 0.0153. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

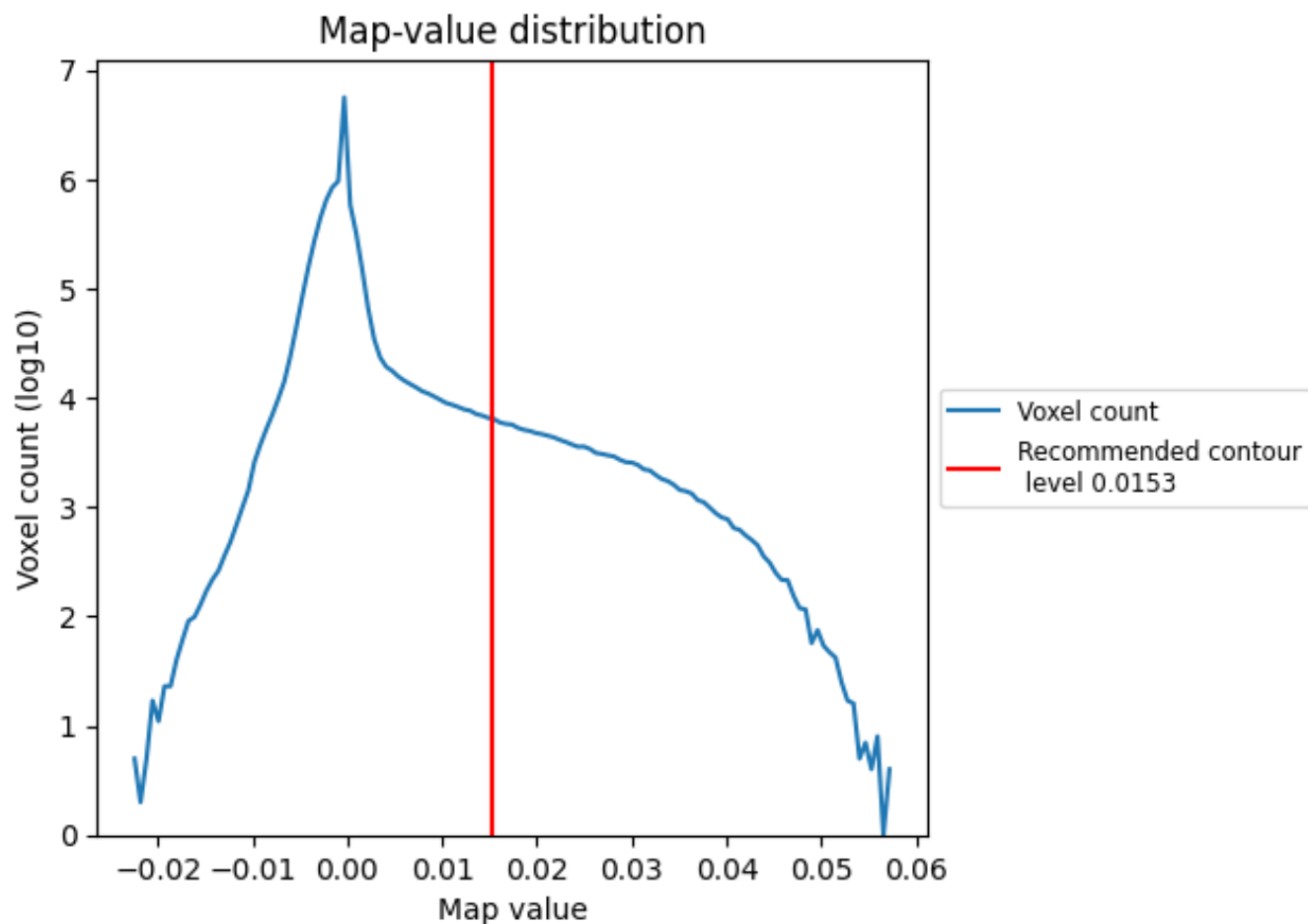
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

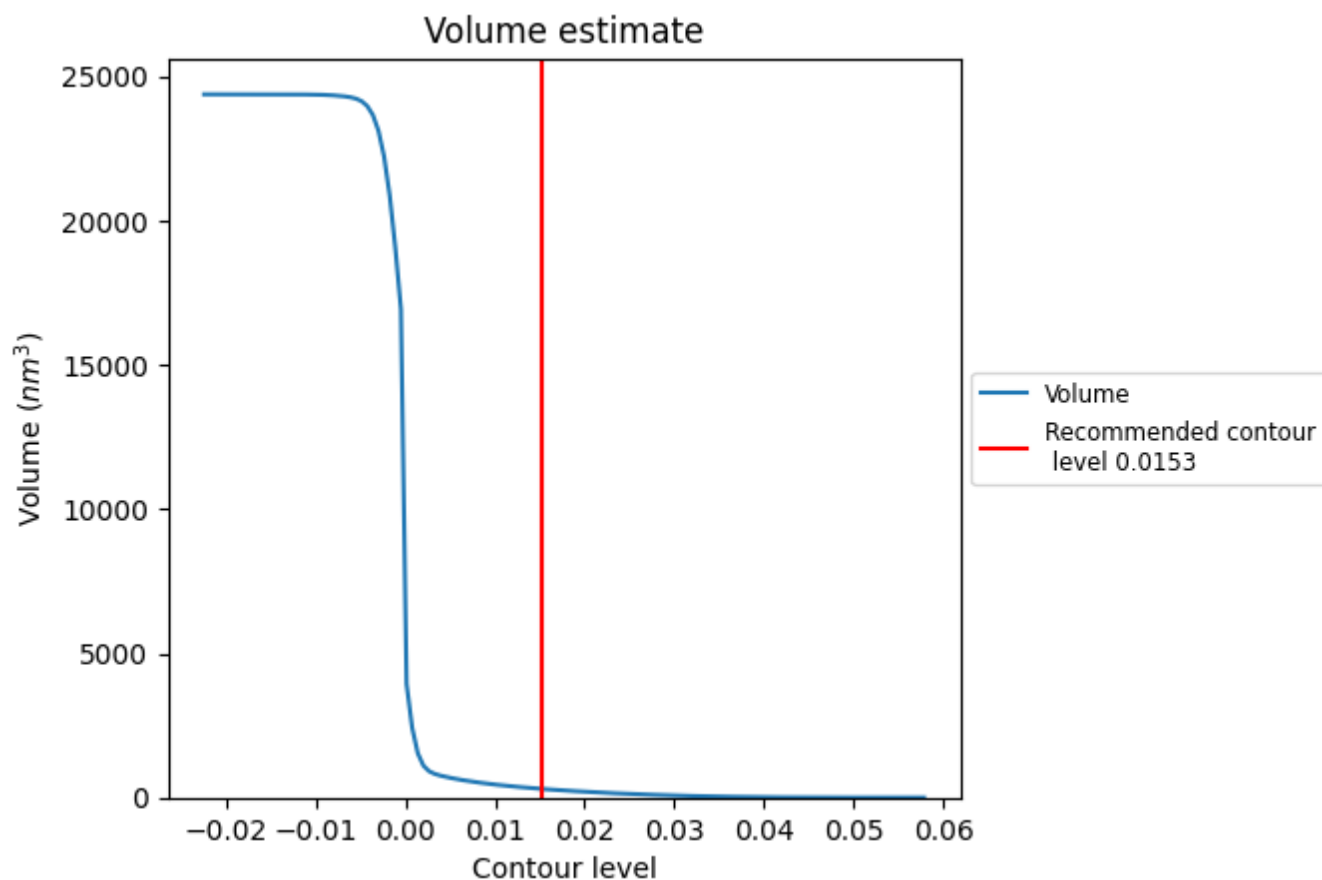
This section contains the results of statistical analysis of the map.

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.

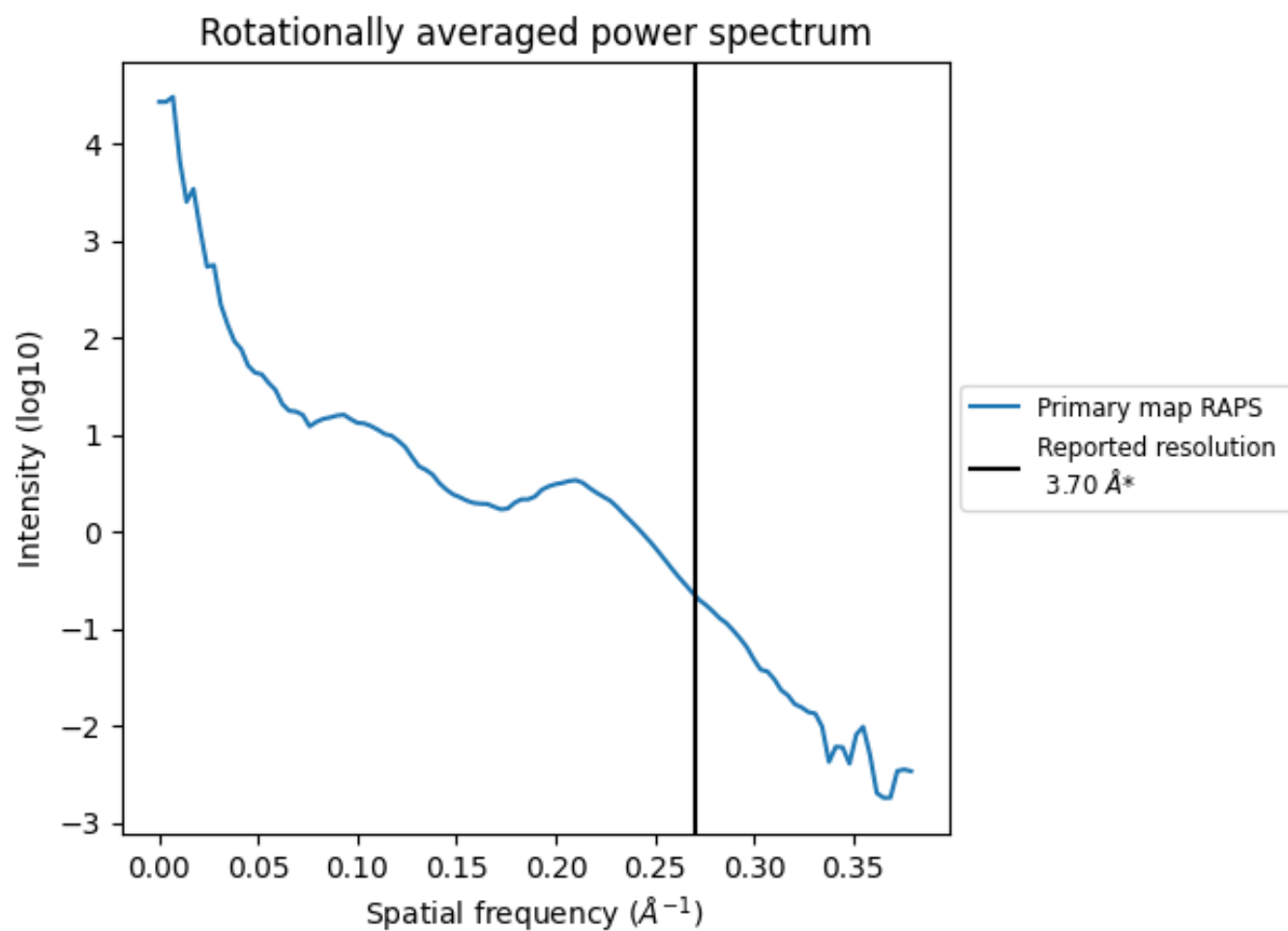
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 298 nm³; this corresponds to an approximate mass of 270 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 0.270 Å⁻¹

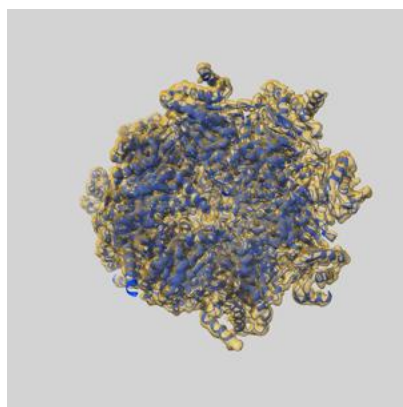
8 Fourier-Shell correlation

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

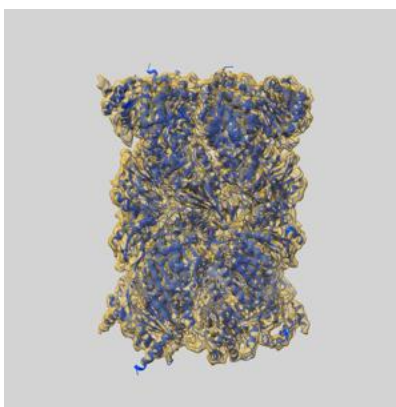
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-13389 and PDB model 7PG9. Per-residue inclusion information can be found in section [3](#) on page [7](#).

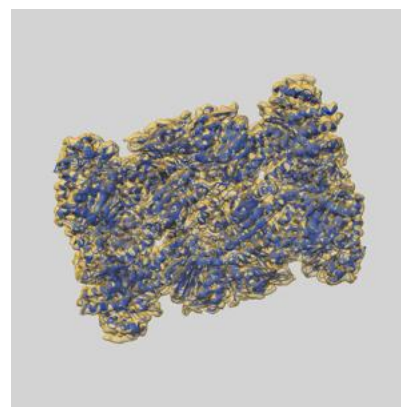
9.1 Map-model overlay [i](#)



X



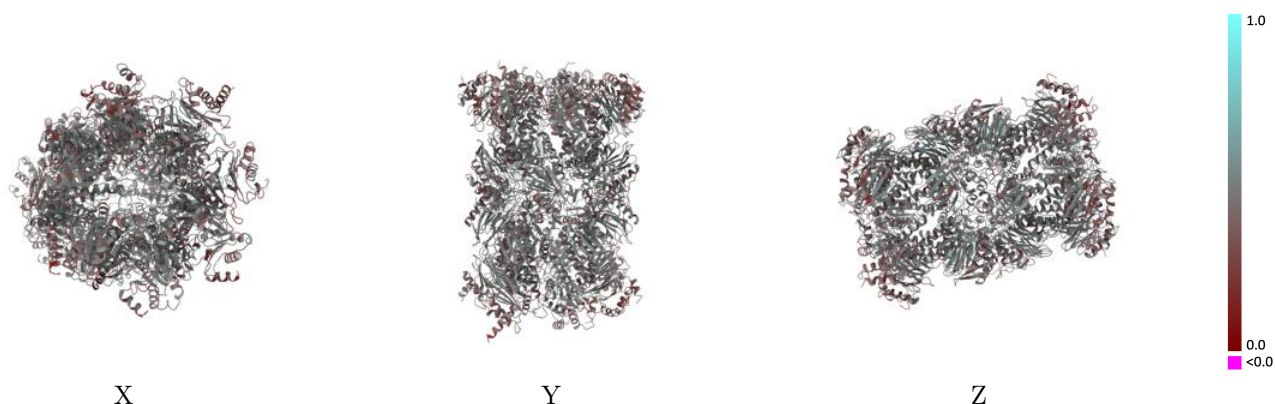
Y



Z

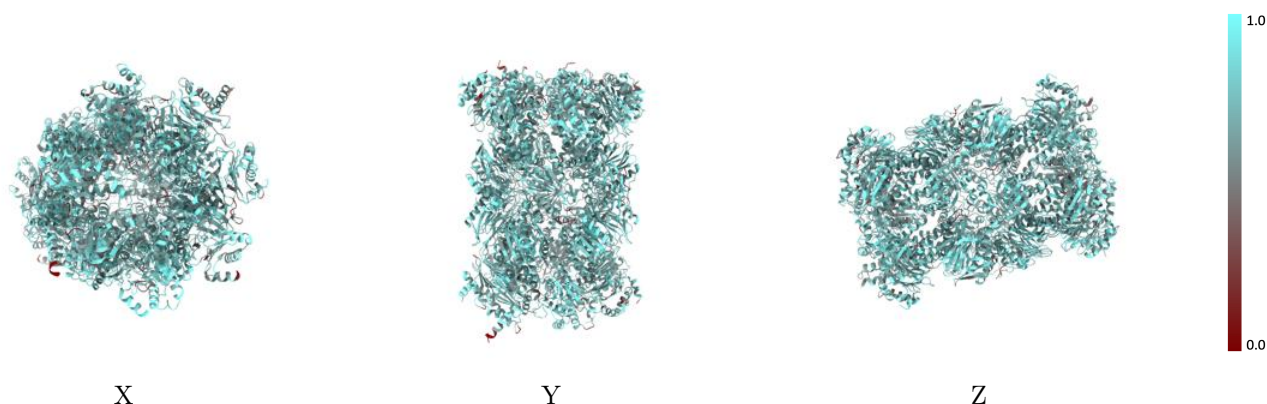
The images above show the 3D surface view of the map at the recommended contour level 0.0153 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



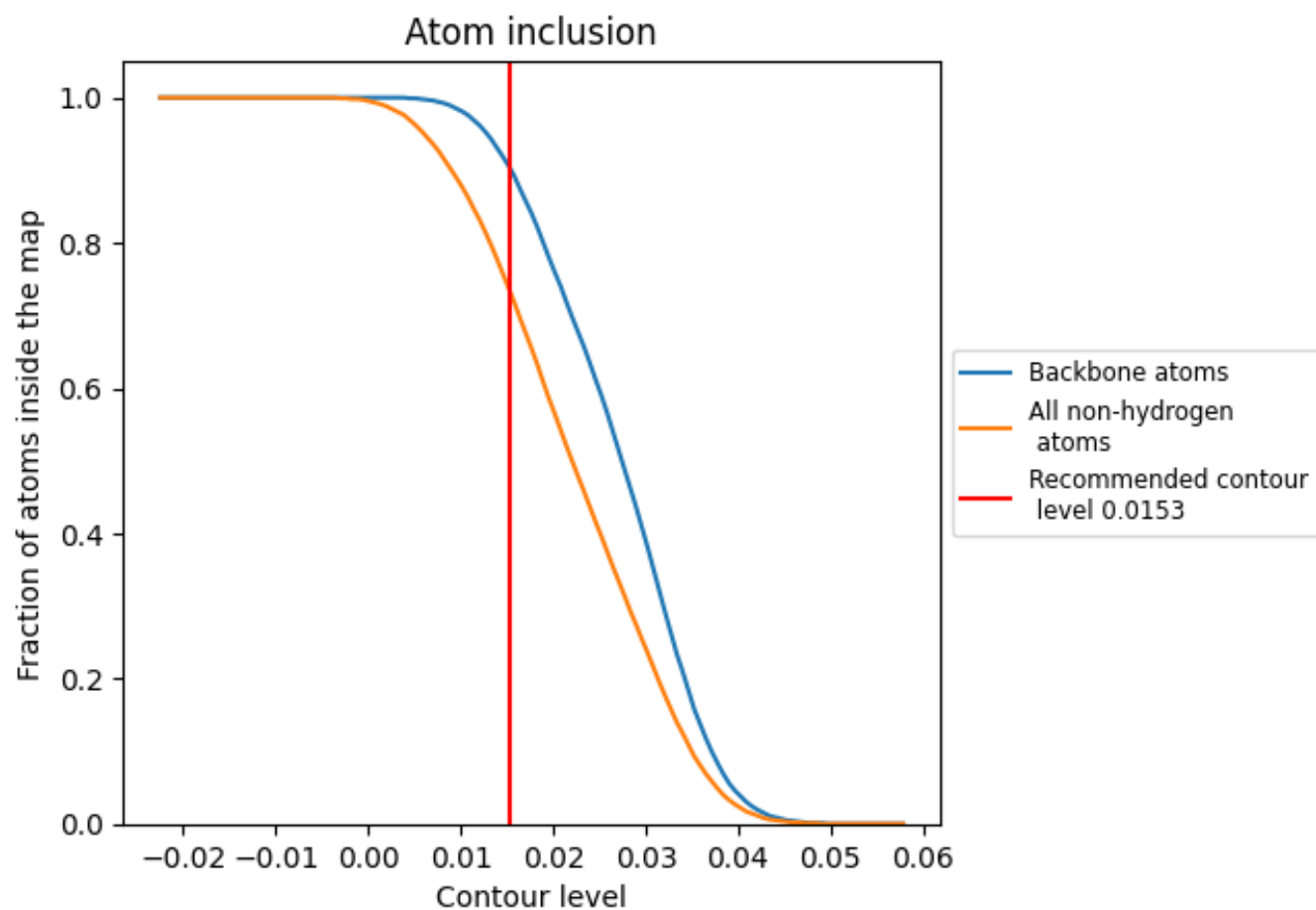
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0153).



























































9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 74% 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 (0.0153) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7350	 0.4440
A	 0.6860	 0.4190
B	 0.7430	 0.4480
C	 0.6880	 0.4330
D	 0.7340	 0.4470
E	 0.7000	 0.4250
F	 0.7290	 0.4300
G	 0.7210	 0.4300
H	 0.7400	 0.4580
I	 0.7380	 0.4510
J	 0.7380	 0.4560
K	 0.7390	 0.4550
L	 0.7710	 0.4600
M	 0.7240	 0.4500
N	 0.7600	 0.4620
O	 0.7110	 0.4300
P	 0.7590	 0.4500
Q	 0.7110	 0.4290
R	 0.7570	 0.4430
S	 0.7100	 0.4300
T	 0.7510	 0.4350
U	 0.7500	 0.4330
V	 0.7410	 0.4490
W	 0.7390	 0.4500
X	 0.7400	 0.4540
Y	 0.7530	 0.4600
Z	 0.7780	 0.4660
a	 0.7280	 0.4460
b	 0.7710	 0.4580

