



wwPDB EM Validation Summary Report ⓘ

Nov 6, 2022 – 11:07 PM EST

PDB ID : 6MUX
EMDB ID : EMD-9259
Title : The structure of the Plasmodium falciparum 20S proteasome in complex with one PA28 activator
Authors : Metcalfe, R.D.; Xie, S.C.; Hanssen, E.; Gillett, D.L.; Leis, A.P.; Tilley, L.; Griffin, M.D.W.
Deposited on : 2018-10-23
Resolution : 3.90 Å (reported)
Based on initial models : 6DFK, 6MUW

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.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

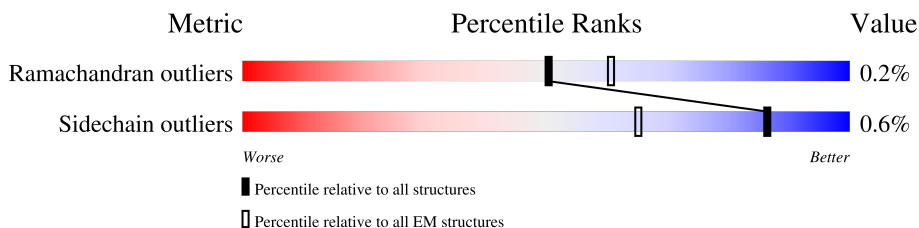
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.90 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 ≥ 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	260	
1	O	260	
2	B	235	
2	P	235	
3	C	246	
3	Q	246	
4	D	241	
4	R	241	
5	E	256	

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Mol	Chain	Length	Quality of chain
5	S	256	
6	F	254	
6	T	254	
7	G	252	
7	U	252	
8	H	252	
8	V	252	
9	I	229	
9	W	229	
10	J	218	
10	X	218	
11	K	195	
11	Y	195	
12	L	211	
12	Z	211	
13	M	240	
13	a	240	
14	N	265	
14	b	265	
15	c	279	
15	d	279	
15	e	279	
15	f	279	
15	g	279	
15	h	279	

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Mol	Chain	Length	Quality of chain
15	i	279	<div><div><div></div><div></div><div></div></div><div>42%67%.</div><div>32%</div></div>

2 Entry composition

There are 15 unique types of molecules in this entry. The entry contains 60284 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 20S proteasome alpha-1 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	244	Total	C	N	O	S	0	0
			1928	1213	323	377	15		
1	O	244	Total	C	N	O	S	0	0
			1928	1213	323	377	15		

- Molecule 2 is a protein called 20S proteasome alpha-2 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	224	Total	C	N	O	S	0	0
			1782	1146	293	337	6		
2	P	229	Total	C	N	O	S	0	0
			1826	1175	298	347	6		

- Molecule 3 is a protein called 20S proteasome alpha-3 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	240	Total	C	N	O	S	0	0
			1918	1230	308	377	3		
3	Q	245	Total	C	N	O	S	0	0
			1965	1258	319	385	3		

- Molecule 4 is a protein called 20S proteasome alpha-4 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	227	Total	C	N	O	S	0	0
			1795	1147	303	337	8		
4	R	233	Total	C	N	O	S	0	0
			1845	1178	312	347	8		

- Molecule 5 is a protein called 20S proteasome alpha-5 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	227	Total	C	N	O	S	0	0
			1760	1109	291	350	10		
5	S	228	Total	C	N	O	S	0	0
			1767	1115	292	349	11		

- Molecule 6 is a protein called 20S proteasome alpha-6 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	235	Total	C	N	O	S	0	0
			1870	1191	308	361	10		
6	T	234	Total	C	N	O	S	0	0
			1858	1182	307	359	10		

- Molecule 7 is a protein called 20S proteasome alpha-7 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	233	Total	C	N	O	S	0	0
			1906	1215	319	360	12		
7	U	233	Total	C	N	O	S	0	0
			1906	1215	319	360	12		

- Molecule 8 is a protein called 20S proteasome beta-1 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	194	Total	C	N	O	S	0	0
			1547	978	265	292	12		
8	V	194	Total	C	N	O	S	0	0
			1547	978	265	292	12		

- Molecule 9 is a protein called 20S proteasome beta-2 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	219	Total	C	N	O	S	0	0
			1676	1058	291	313	14		
9	W	219	Total	C	N	O	S	0	0
			1676	1058	291	313	14		

- Molecule 10 is a protein called 20S proteasome beta-3 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	203	Total	C	N	O	S	0	0
			1595	1017	258	306	14		

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Mol	Chain	Residues	Atoms					AltConf	Trace
10	X	203	Total	C	N	O	S	0	0
			1595	1017	258	306	14		

- Molecule 11 is a protein called 20S proteasome beta-4 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	195	Total	C	N	O	S	0	0
			1614	1042	266	298	8		
11	Y	195	Total	C	N	O	S	0	0
			1614	1042	266	298	8		

- Molecule 12 is a protein called 20S proteasome beta-5 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	204	Total	C	N	O	S	0	0
			1600	1021	265	307	7		
12	Z	204	Total	C	N	O	S	0	0
			1600	1021	265	307	7		

- Molecule 13 is a protein called 20S proteasome beta-6 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	211	Total	C	N	O	S	0	0
			1676	1073	277	319	7		
13	a	211	Total	C	N	O	S	0	0
			1676	1073	277	319	7		

- Molecule 14 is a protein called 20S proteasome beta-7 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	224	Total	C	N	O	S	0	0
			1845	1177	313	348	7		
14	b	224	Total	C	N	O	S	0	0
			1845	1177	313	348	7		

- Molecule 15 is a protein called Proteasome activator PA28.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	c	188	Total	C	N	O	S	0	0
			1578	1019	261	295	3		
15	d	190	Total	C	N	O	S	0	0
			1595	1029	265	298	3		

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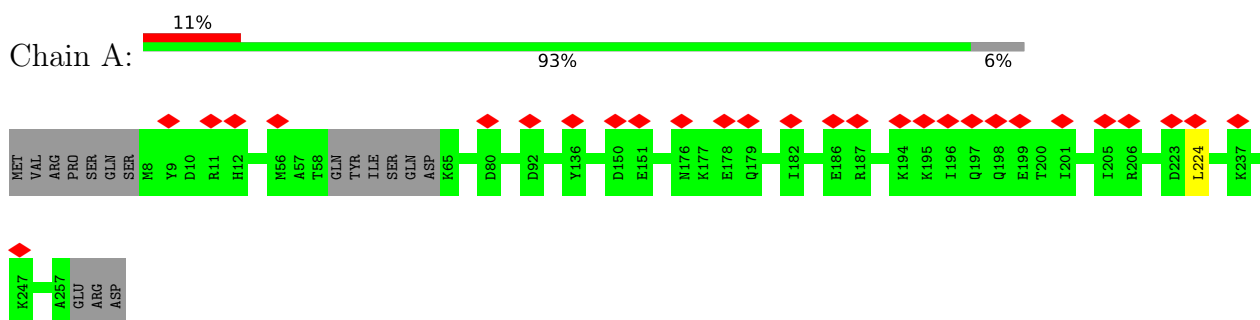
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Mol	Chain	Residues	Atoms					AltConf	Trace
15	e	187	Total 1571	C 1014	N 260	O 294	S 3	0	0
15	f	190	Total 1595	C 1029	N 265	O 298	S 3	0	0
15	g	190	Total 1595	C 1029	N 265	O 298	S 3	0	0
15	h	190	Total 1595	C 1029	N 265	O 298	S 3	0	0
15	i	190	Total 1595	C 1029	N 265	O 298	S 3	0	0

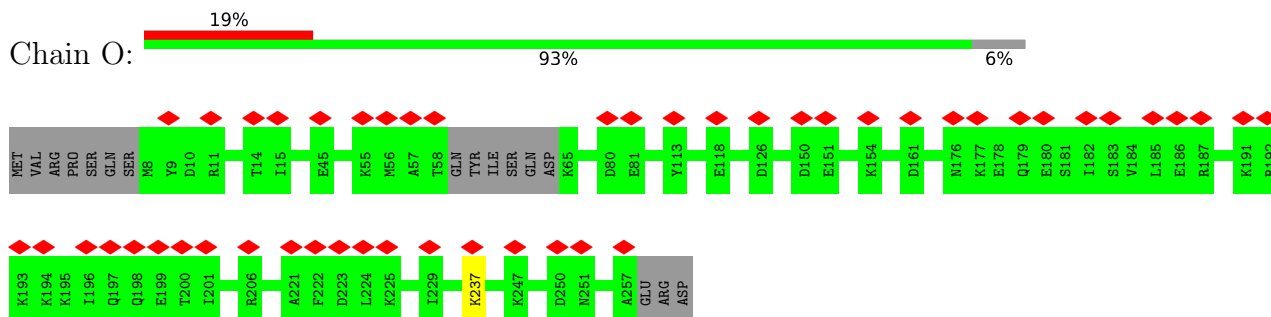
3 Residue-property plots [i](#)

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

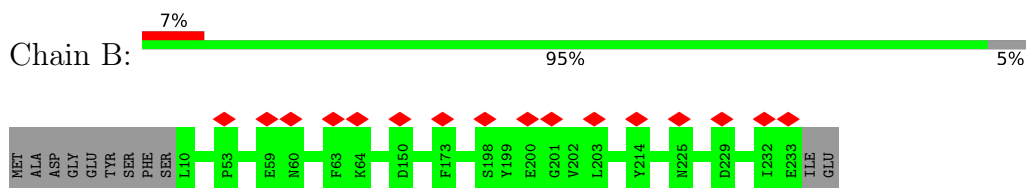
- Molecule 1: 20S proteasome alpha-1 subunit



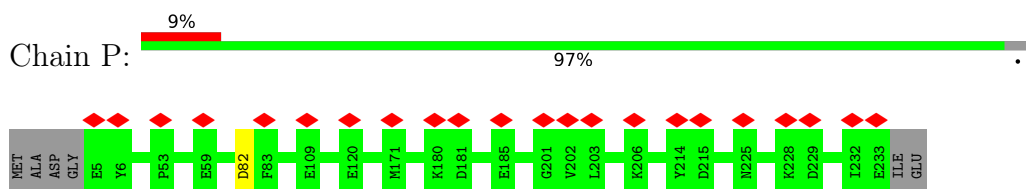
- Molecule 1: 20S proteasome alpha-1 subunit



- Molecule 2: 20S proteasome alpha-2 subunit

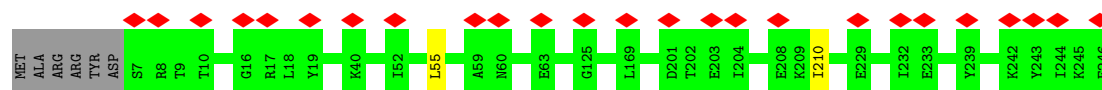


- Molecule 2: 20S proteasome alpha-2 subunit



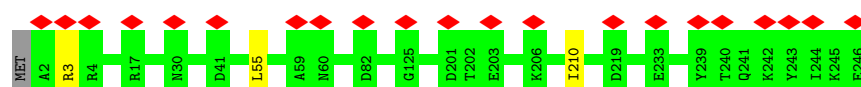
- Molecule 3: 20S proteasome alpha-3 subunit

Chain C: 

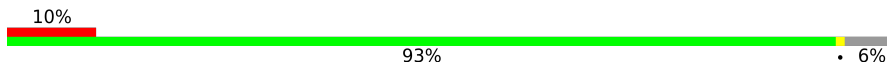


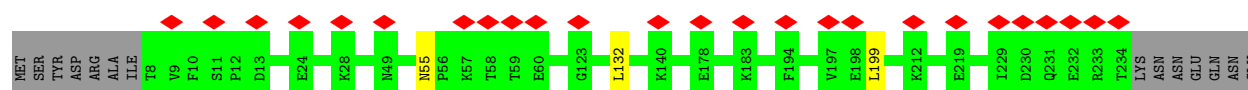
- Molecule 3: 20S proteasome alpha-3 subunit

Chain Q: 



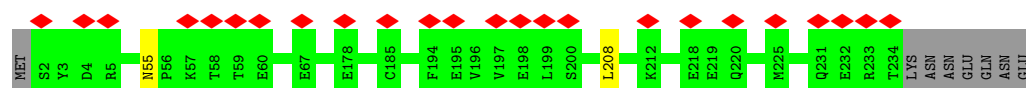
- Molecule 4: 20S proteasome alpha-4 subunit

Chain D: 




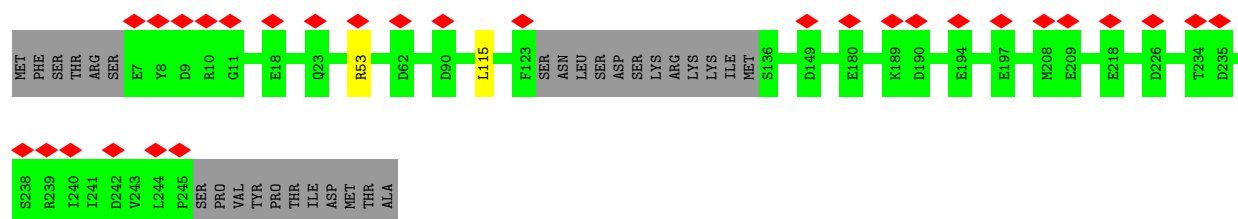
- Molecule 4: 20S proteasome alpha-4 subunit

Chain R: 




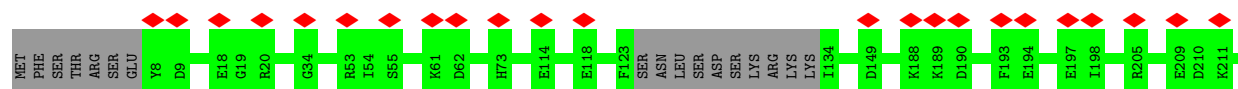
- Molecule 5: 20S proteasome alpha-5 subunit

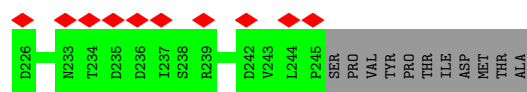
Chain E: 



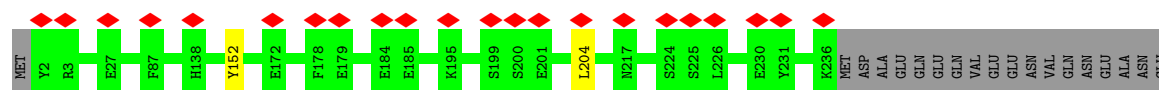
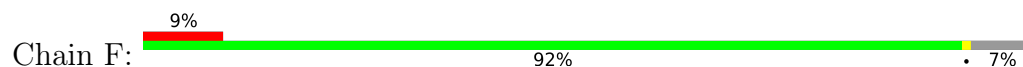
- Molecule 5: 20S proteasome alpha-5 subunit

Chain S: 

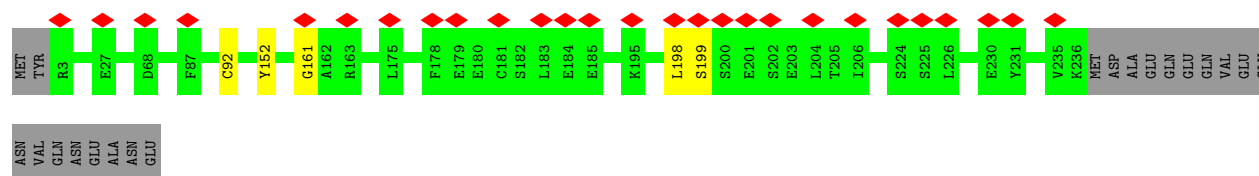




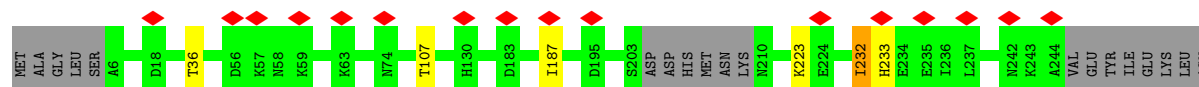
- Molecule 6: 20S proteasome alpha-6 subunit



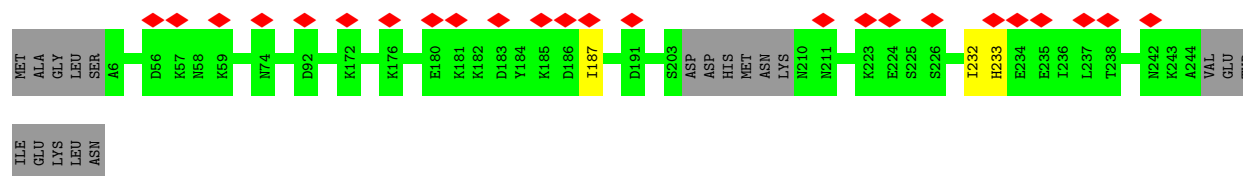
- Molecule 6: 20S proteasome alpha-6 subunit



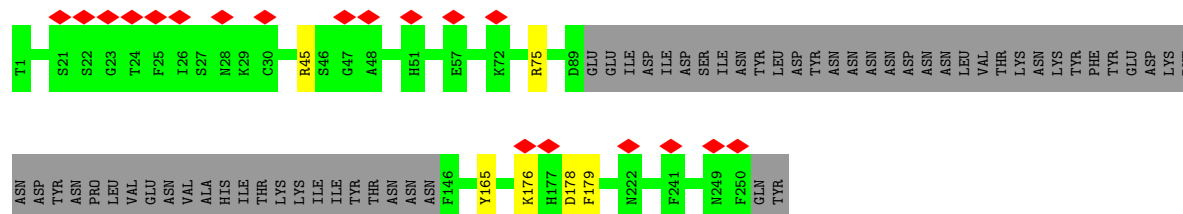
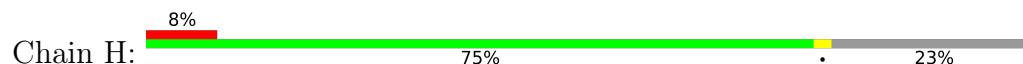
- Molecule 7: 20S proteasome alpha-7 subunit



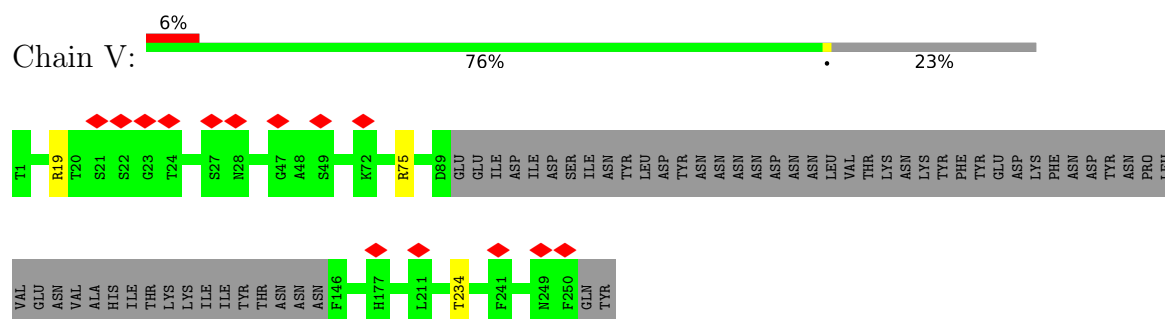
- Molecule 7: 20S proteasome alpha-7 subunit



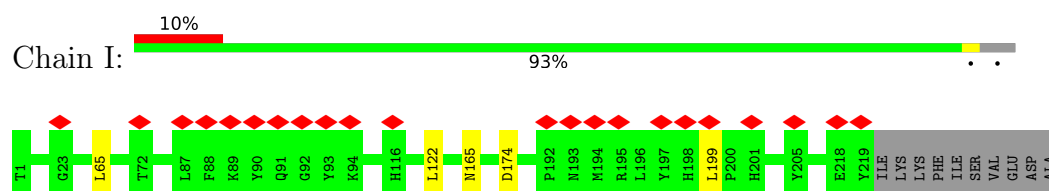
- Molecule 8: 20S proteasome beta-1 subunit



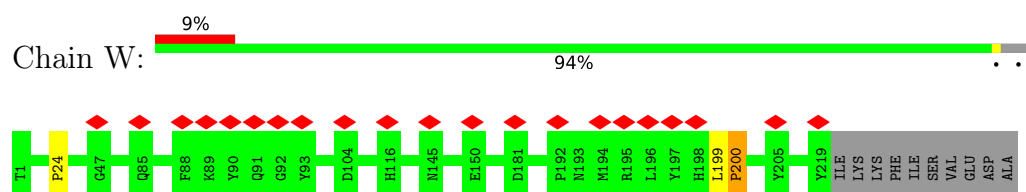
- Molecule 8: 20S proteasome beta-1 subunit



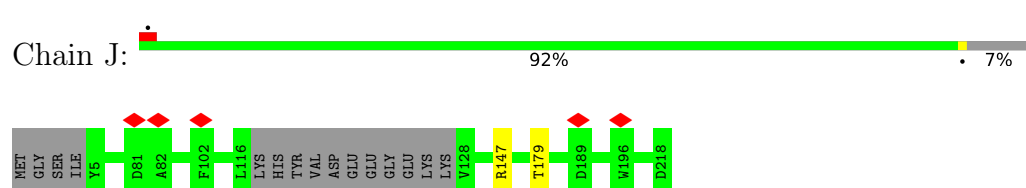
- Molecule 9: 20S proteasome beta-2 subunit



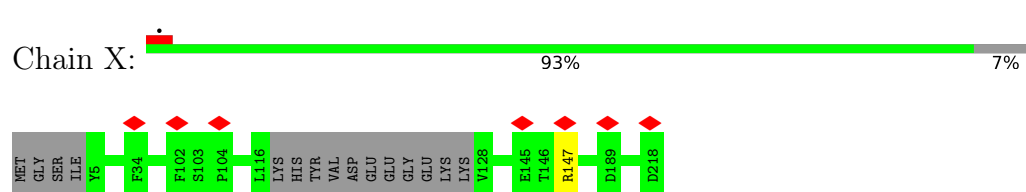
- Molecule 9: 20S proteasome beta-2 subunit



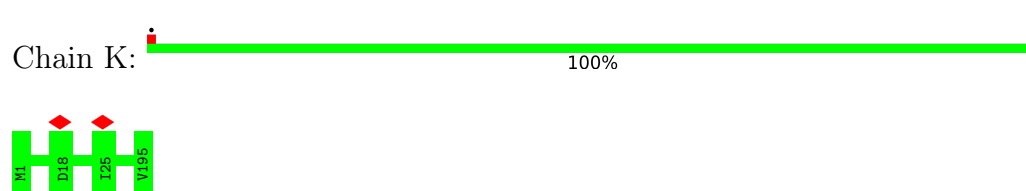
- Molecule 10: 20S proteasome beta-3 subunit



- Molecule 10: 20S proteasome beta-3 subunit

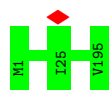


- Molecule 11: 20S proteasome beta-4 subunit



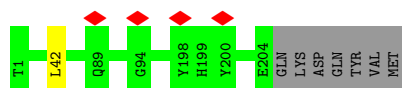
- Molecule 11: 20S proteasome beta-4 subunit

Chain Y:  100%



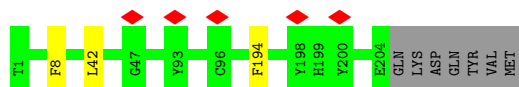
- Molecule 12: 20S proteasome beta-5 subunit

Chain L:  96%



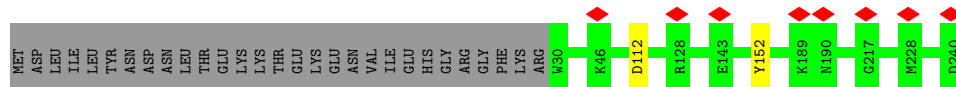
- Molecule 12: 20S proteasome beta-5 subunit

Chain Z:  95%


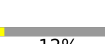


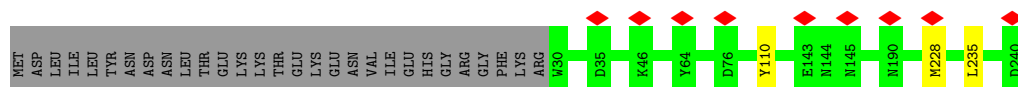
- Molecule 13: 20S proteasome beta-6 subunit

Chain M:  87%  12%



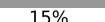


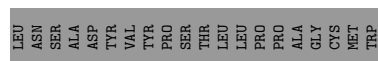
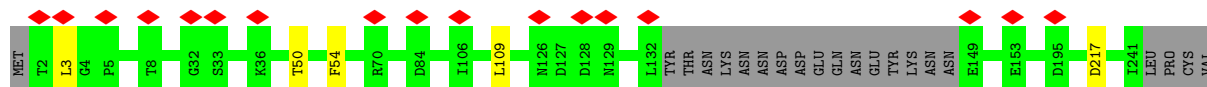
- Molecule 13: 20S proteasome beta-6 subunit

Chain a:  87%  12%



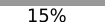


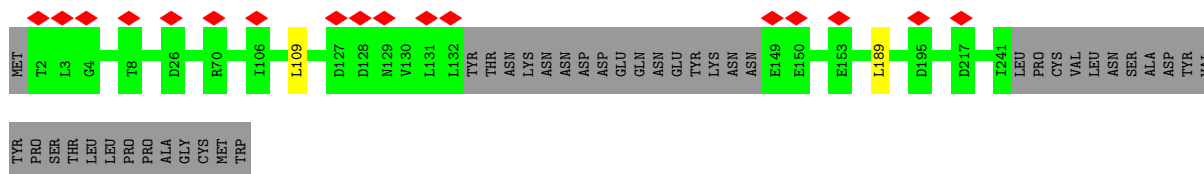
- Molecule 14: 20S proteasome beta-7 subunit

Chain N:  6%  83%  15%

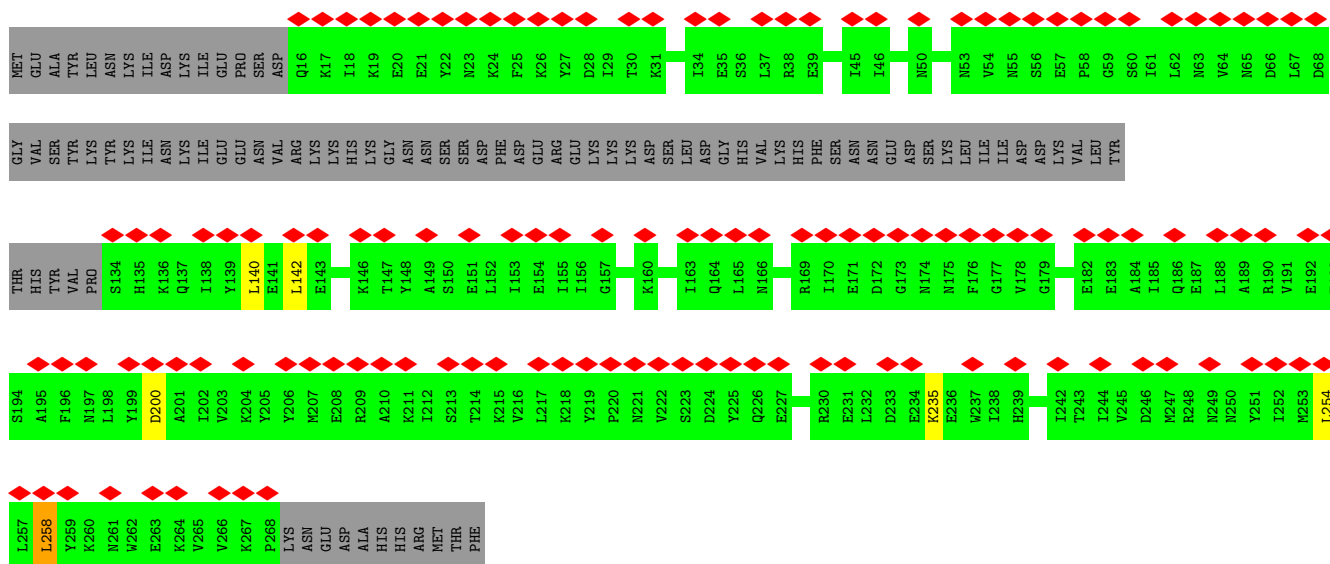


- Molecule 14: 20S proteasome beta-7 subunit

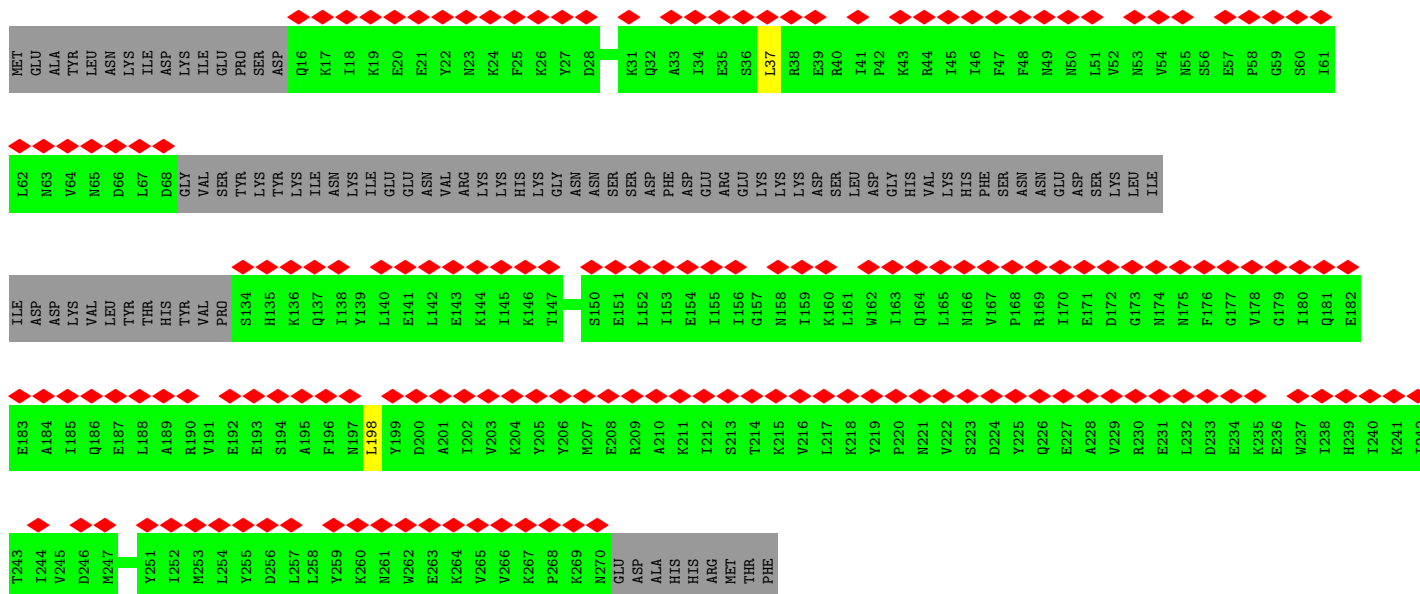
Chain b:  6%  84%  15%



• Molecule 15: Proteasome activator PA28

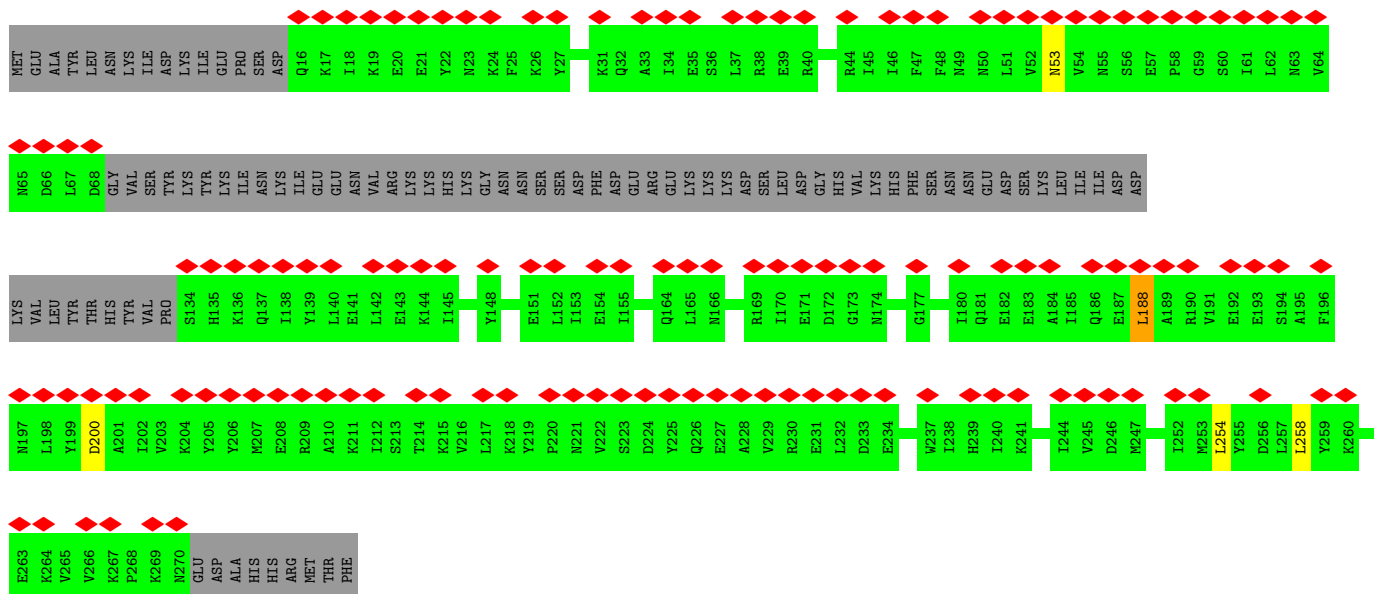


• Molecule 15: Proteasome activator PA28

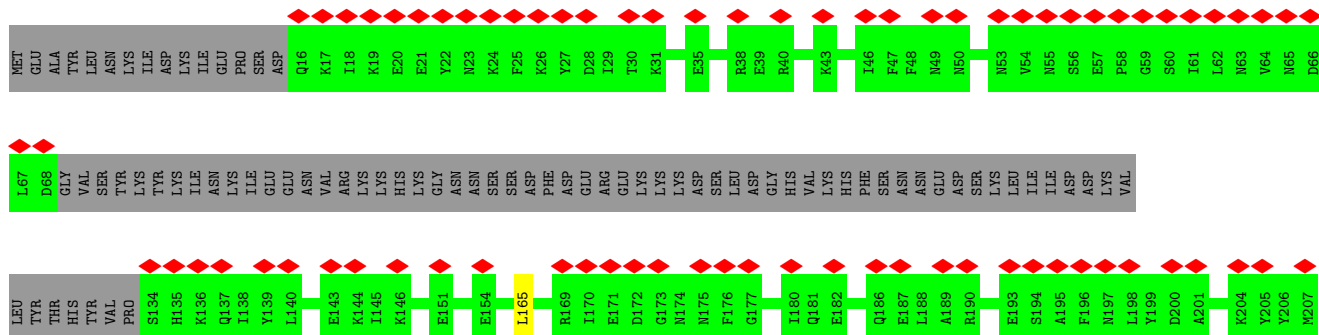
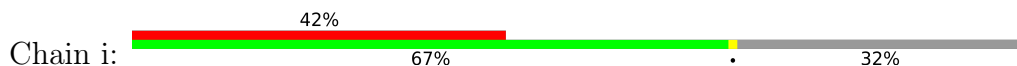


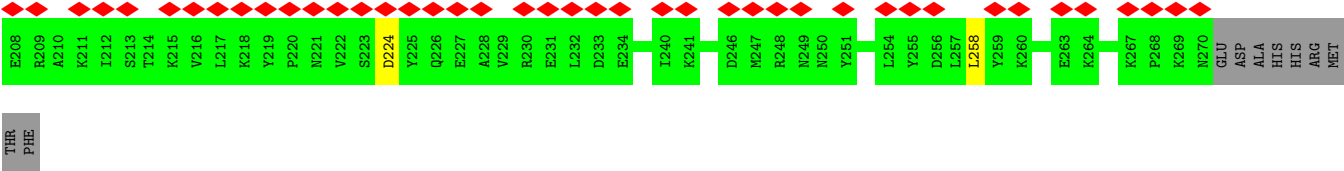
• Molecule 15: Proteasome activator PA28

- Molecule 15: Proteasome activator PA28



- Molecule 15: Proteasome activator PA28





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	57337	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	32	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	100000	Depositor
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	33.420	Depositor
Minimum map value	-19.000	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	6.0	Depositor
Map size (\AA)	450.63998, 450.63998, 450.63998	wwPDB
Map dimensions	344, 344, 344	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.31, 1.31, 1.31	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	0.41	0/1953	0.64	1/2632 (0.0%)
1	O	0.38	0/1953	0.66	0/2632
2	B	0.43	0/1814	0.63	0/2450
2	P	0.41	0/1860	0.63	0/2512
3	C	0.48	0/1953	0.64	1/2643 (0.0%)
3	Q	0.45	0/2001	0.65	1/2707 (0.0%)
4	D	0.42	0/1824	0.64	2/2461 (0.1%)
4	R	0.40	0/1875	0.63	1/2530 (0.0%)
5	E	0.37	0/1784	0.60	1/2410 (0.0%)
5	S	0.37	0/1791	0.59	0/2419
6	F	0.48	1/1905 (0.1%)	0.64	0/2566
6	T	0.45	1/1892 (0.1%)	0.65	2/2548 (0.1%)
7	G	0.46	0/1947	0.71	1/2631 (0.0%)
7	U	0.44	0/1947	0.67	0/2631
8	H	0.50	0/1571	0.69	1/2105 (0.0%)
8	V	0.45	0/1571	0.66	0/2105
9	I	0.42	1/1712 (0.1%)	0.67	2/2328 (0.1%)
9	W	0.41	1/1712 (0.1%)	0.65	1/2328 (0.0%)
10	J	0.53	0/1621	0.67	0/2189
10	X	0.53	0/1621	0.66	0/2189
11	K	0.55	0/1649	0.67	0/2223
11	Y	0.54	0/1649	0.67	0/2223
12	L	0.50	0/1633	0.68	1/2202 (0.0%)
12	Z	0.55	1/1633 (0.1%)	0.69	1/2202 (0.0%)
13	M	0.47	0/1708	0.67	0/2314
13	a	0.49	0/1708	0.69	0/2314
14	N	0.44	0/1882	0.69	4/2538 (0.2%)
14	b	0.44	0/1882	0.64	2/2538 (0.1%)
15	c	0.37	0/1606	0.69	4/2166 (0.2%)
15	d	0.34	0/1623	0.67	2/2188 (0.1%)
15	e	0.33	0/1598	0.68	1/2154 (0.0%)
15	f	0.32	0/1623	0.66	2/2188 (0.1%)
15	g	0.33	0/1623	0.67	4/2188 (0.2%)
15	h	0.35	0/1623	0.65	3/2188 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
15	i	0.35	0/1623	0.65	2/2188 (0.1%)
All	All	0.44	5/61370 (0.0%)	0.66	40/82830 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	T	152	TYR	CD1-CE1	-6.05	1.30	1.39
9	I	199	LEU	C-N	5.70	1.45	1.34
6	F	152	TYR	CD1-CE1	-5.44	1.31	1.39
9	W	199	LEU	C-N	5.44	1.44	1.34
12	Z	8	PHE	CB-CG	-5.16	1.42	1.51

The worst 5 of 40 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	N	109	LEU	CA-CB-CG	9.67	137.54	115.30
14	b	109	LEU	CA-CB-CG	9.00	136.00	115.30
15	f	254	LEU	CA-CB-CG	8.37	134.54	115.30
15	i	165	LEU	CA-CB-CG	7.62	132.84	115.30
3	Q	55	LEU	CA-CB-CG	7.16	131.76	115.30

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	240/260 (92%)	214 (89%)	26 (11%)	0	100	100
1	O	240/260 (92%)	211 (88%)	29 (12%)	0	100	100
2	B	222/235 (94%)	203 (91%)	19 (9%)	0	100	100
2	P	227/235 (97%)	205 (90%)	22 (10%)	0	100	100
3	C	238/246 (97%)	217 (91%)	20 (8%)	1 (0%)	34	71
3	Q	243/246 (99%)	218 (90%)	24 (10%)	1 (0%)	34	71
4	D	225/241 (93%)	200 (89%)	25 (11%)	0	100	100
4	R	231/241 (96%)	211 (91%)	20 (9%)	0	100	100
5	E	223/256 (87%)	197 (88%)	26 (12%)	0	100	100
5	S	224/256 (88%)	200 (89%)	24 (11%)	0	100	100
6	F	233/254 (92%)	195 (84%)	37 (16%)	1 (0%)	34	71
6	T	232/254 (91%)	204 (88%)	27 (12%)	1 (0%)	34	71
7	G	229/252 (91%)	189 (82%)	37 (16%)	3 (1%)	12	48
7	U	229/252 (91%)	197 (86%)	29 (13%)	3 (1%)	12	48
8	H	190/252 (75%)	159 (84%)	29 (15%)	2 (1%)	14	51
8	V	190/252 (75%)	162 (85%)	28 (15%)	0	100	100
9	I	217/229 (95%)	175 (81%)	42 (19%)	0	100	100
9	W	217/229 (95%)	190 (88%)	26 (12%)	1 (0%)	29	67
10	J	199/218 (91%)	175 (88%)	24 (12%)	0	100	100
10	X	199/218 (91%)	176 (88%)	23 (12%)	0	100	100
11	K	193/195 (99%)	171 (89%)	22 (11%)	0	100	100
11	Y	193/195 (99%)	168 (87%)	25 (13%)	0	100	100
12	L	202/211 (96%)	166 (82%)	36 (18%)	0	100	100
12	Z	202/211 (96%)	171 (85%)	31 (15%)	0	100	100
13	M	209/240 (87%)	184 (88%)	25 (12%)	0	100	100
13	a	209/240 (87%)	178 (85%)	31 (15%)	0	100	100
14	N	220/265 (83%)	189 (86%)	31 (14%)	0	100	100
14	b	220/265 (83%)	192 (87%)	28 (13%)	0	100	100
15	c	184/279 (66%)	178 (97%)	6 (3%)	0	100	100
15	d	186/279 (67%)	180 (97%)	6 (3%)	0	100	100
15	e	183/279 (66%)	177 (97%)	6 (3%)	0	100	100
15	f	186/279 (67%)	179 (96%)	7 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
15	g	186/279 (67%)	180 (97%)	6 (3%)	0	100	100
15	h	186/279 (67%)	179 (96%)	7 (4%)	0	100	100
15	i	186/279 (67%)	180 (97%)	6 (3%)	0	100	100
All	All	7393/8661 (85%)	6570 (89%)	810 (11%)	13 (0%)	50	79

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
7	G	233	HIS
8	H	178	ASP
6	T	199	SER
6	F	204	LEU
7	G	232	ILE

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	215/231 (93%)	215 (100%)	0	100	100
1	O	215/231 (93%)	214 (100%)	1 (0%)	88	93
2	B	196/205 (96%)	196 (100%)	0	100	100
2	P	201/205 (98%)	200 (100%)	1 (0%)	88	93
3	C	208/213 (98%)	208 (100%)	0	100	100
3	Q	212/213 (100%)	211 (100%)	1 (0%)	88	93
4	D	194/207 (94%)	193 (100%)	1 (0%)	88	93
4	R	199/207 (96%)	198 (100%)	1 (0%)	88	93
5	E	195/223 (87%)	194 (100%)	1 (0%)	88	93
5	S	196/223 (88%)	196 (100%)	0	100	100
6	F	210/227 (92%)	210 (100%)	0	100	100
6	T	209/227 (92%)	208 (100%)	1 (0%)	88	93
7	G	212/229 (93%)	209 (99%)	3 (1%)	67	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	U	212/229 (93%)	212 (100%)	0	100	100
8	H	174/231 (75%)	171 (98%)	3 (2%)	60	78
8	V	174/231 (75%)	171 (98%)	3 (2%)	60	78
9	I	185/194 (95%)	183 (99%)	2 (1%)	73	84
9	W	185/194 (95%)	184 (100%)	1 (0%)	88	93
10	J	178/191 (93%)	176 (99%)	2 (1%)	73	84
10	X	178/191 (93%)	177 (99%)	1 (1%)	86	91
11	K	174/174 (100%)	174 (100%)	0	100	100
11	Y	174/174 (100%)	174 (100%)	0	100	100
12	L	169/176 (96%)	169 (100%)	0	100	100
12	Z	169/176 (96%)	168 (99%)	1 (1%)	86	91
13	M	189/216 (88%)	187 (99%)	2 (1%)	73	84
13	a	189/216 (88%)	186 (98%)	3 (2%)	62	79
14	N	201/239 (84%)	200 (100%)	1 (0%)	88	93
14	b	201/239 (84%)	201 (100%)	0	100	100
15	c	175/261 (67%)	172 (98%)	3 (2%)	60	78
15	d	177/261 (68%)	177 (100%)	0	100	100
15	e	174/261 (67%)	171 (98%)	3 (2%)	60	78
15	f	177/261 (68%)	174 (98%)	3 (2%)	60	78
15	g	177/261 (68%)	176 (99%)	1 (1%)	86	91
15	h	177/261 (68%)	174 (98%)	3 (2%)	60	78
15	i	177/261 (68%)	176 (99%)	1 (1%)	86	91
All	All	6648/7739 (86%)	6605 (99%)	43 (1%)	86	91

5 of 43 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
13	a	235	LEU
15	f	26	LYS
15	c	140	LEU
15	e	142	LEU
15	f	235	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 101 such sidechains are listed below:

Mol	Chain	Res	Type
6	T	217	ASN
9	W	106	ASN
15	i	50	ASN
7	U	89	ASN
8	V	219	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.

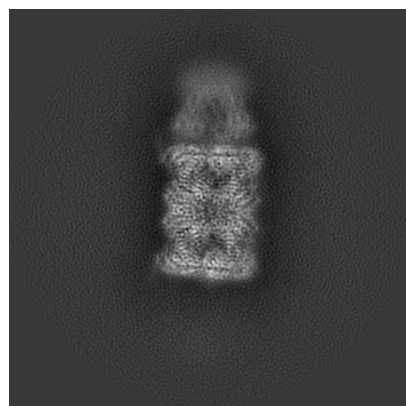
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-9259. These allow visual inspection of the internal detail of the map and identification of artifacts.

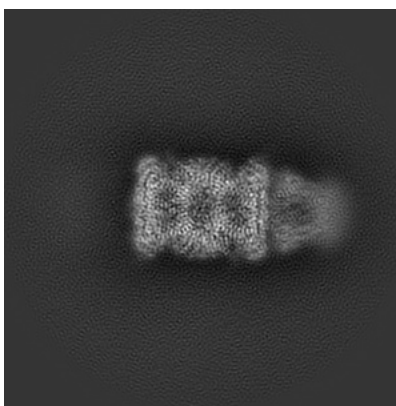
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

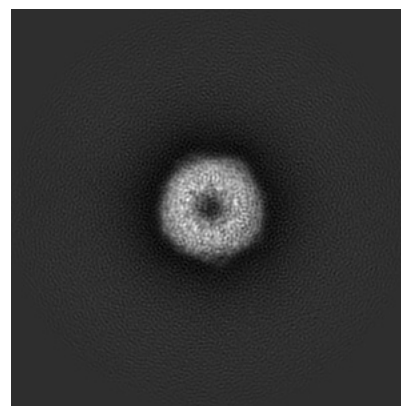
6.1.1 Primary map



X

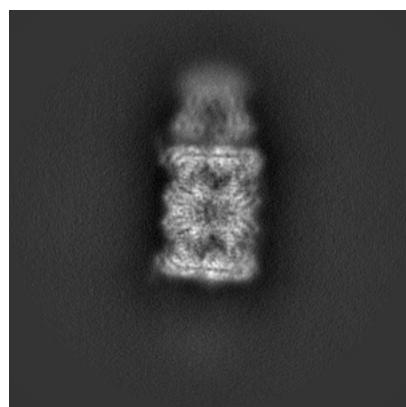


Y

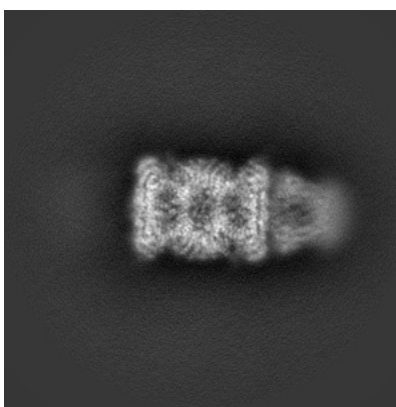


Z

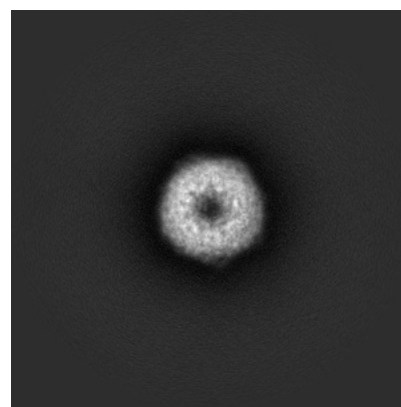
6.1.2 Raw map



X



Y

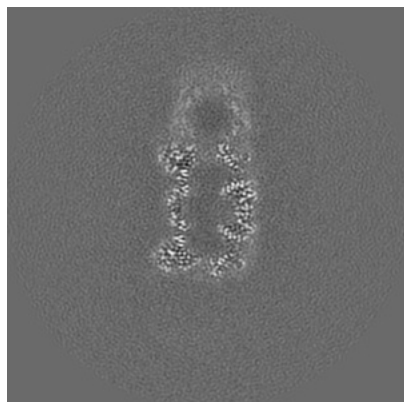


Z

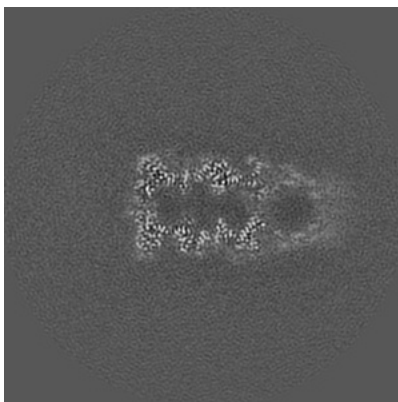
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

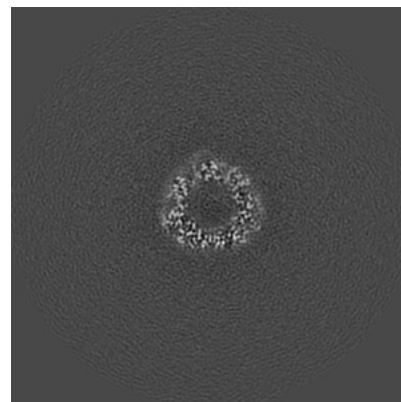
6.2.1 Primary map



X Index: 172

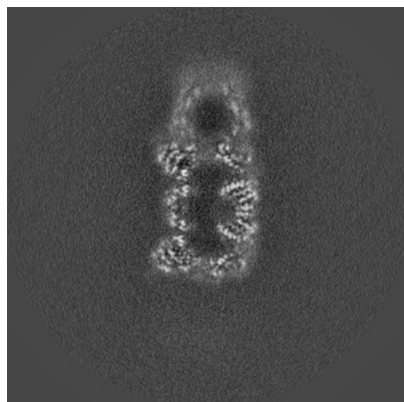


Y Index: 172

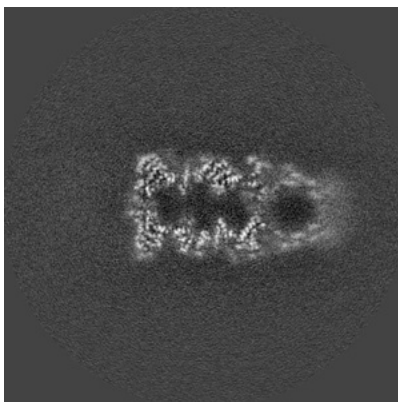


Z Index: 172

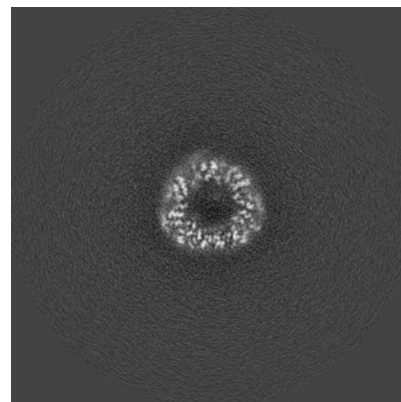
6.2.2 Raw map



X Index: 172



Y Index: 172

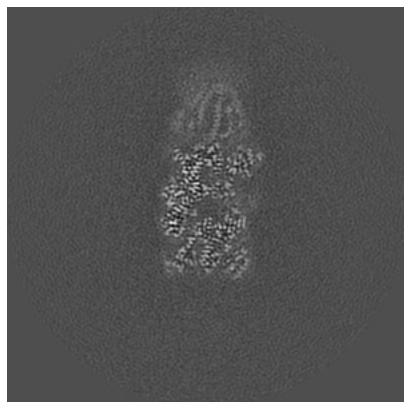


Z Index: 172

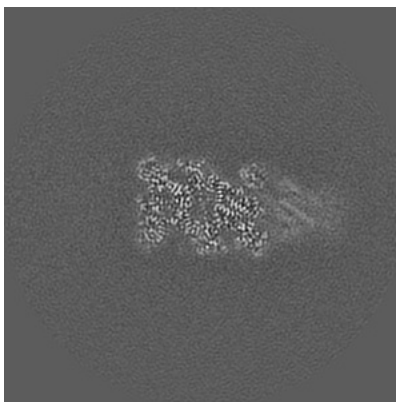
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

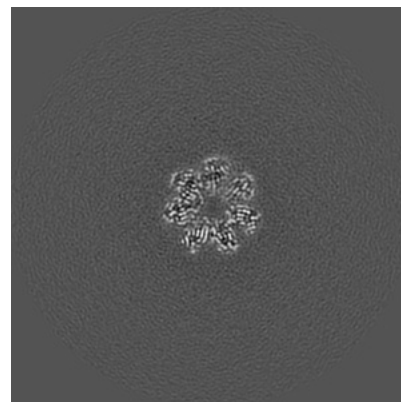
6.3.1 Primary map



X Index: 154

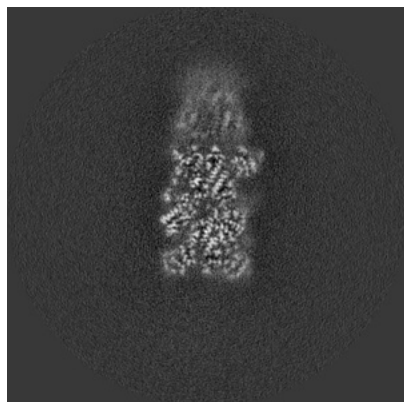


Y Index: 152

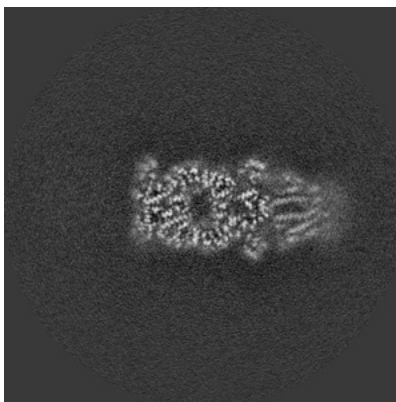


Z Index: 156

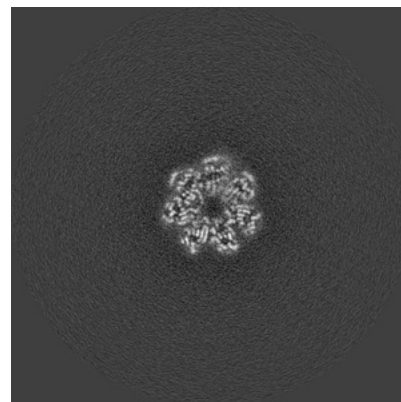
6.3.2 Raw map



X Index: 151



Y Index: 190

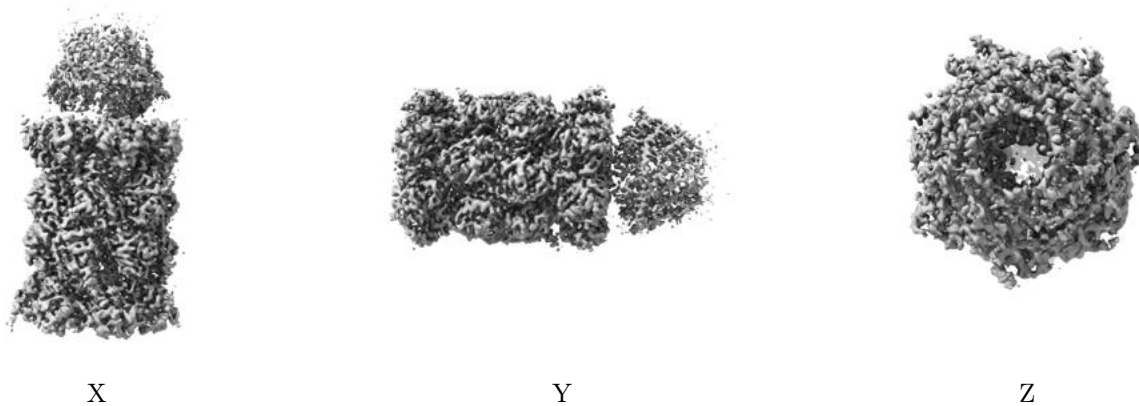


Z Index: 156

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

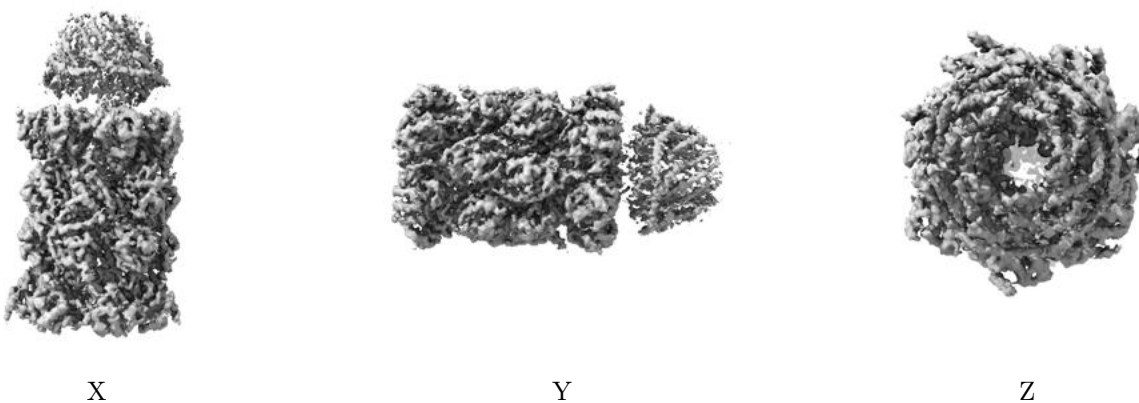
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 6.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

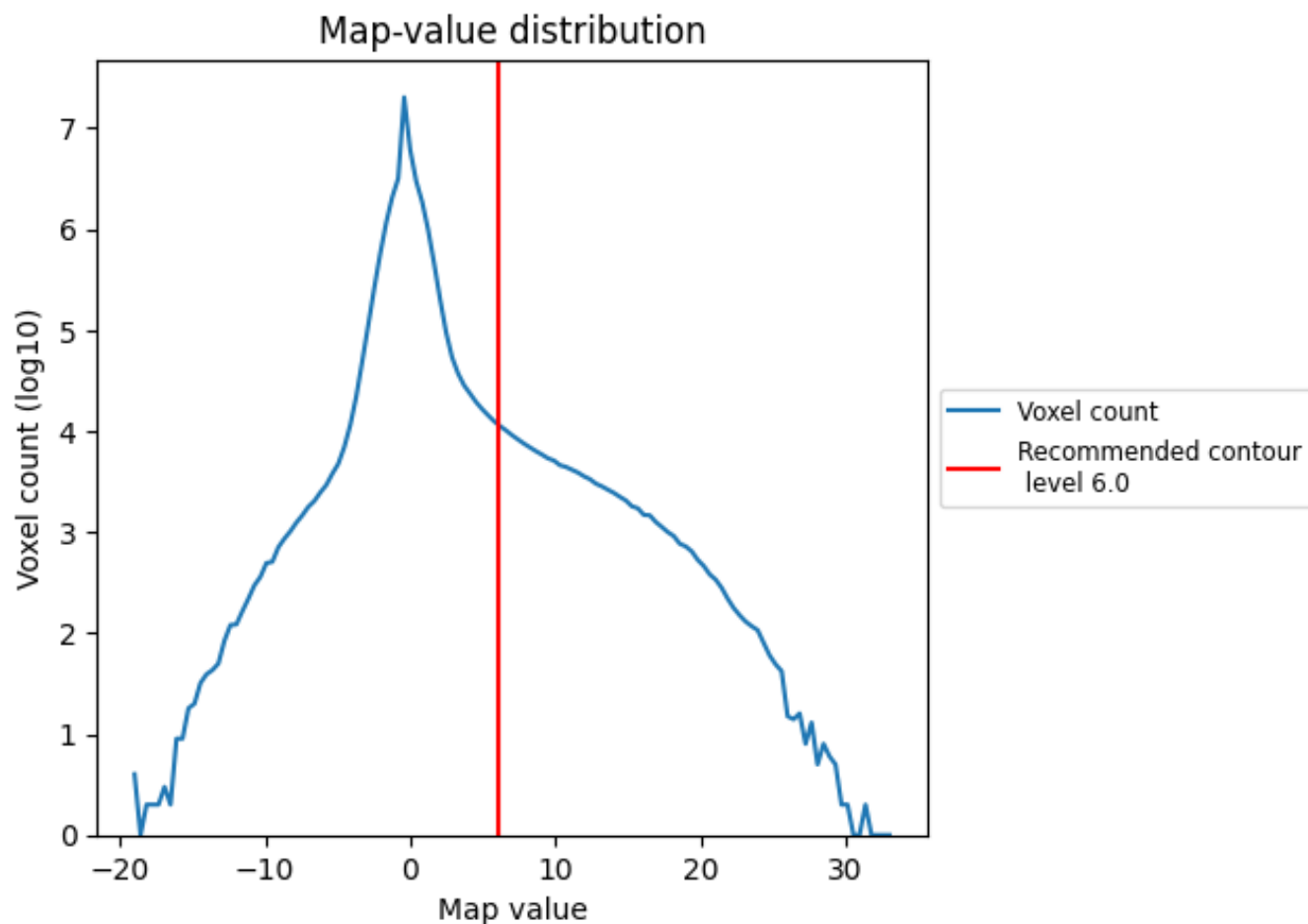
6.5 Mask visualisation [i](#)

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

7 Map analysis [i](#)

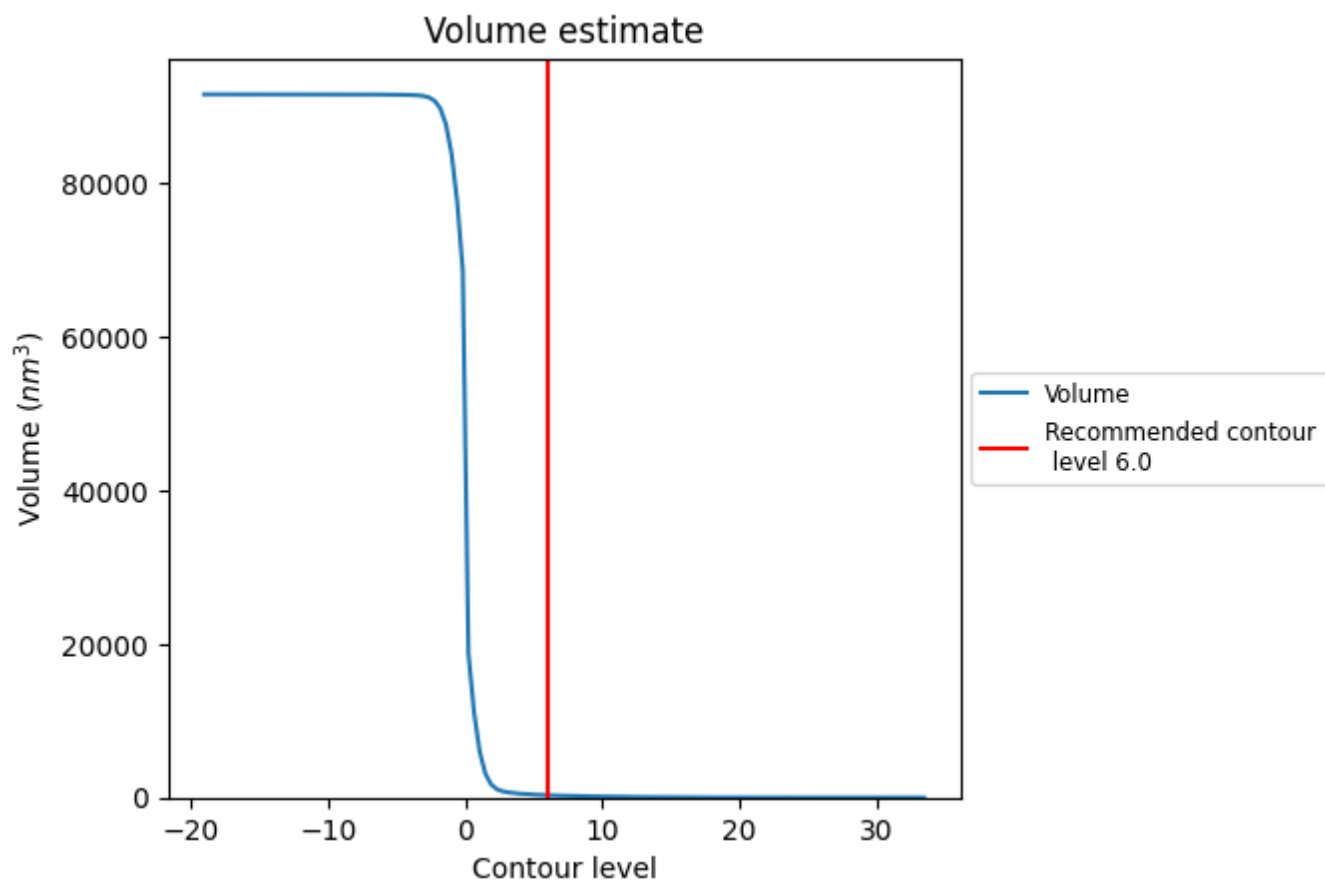
This section contains the results of statistical analysis of the map.

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.

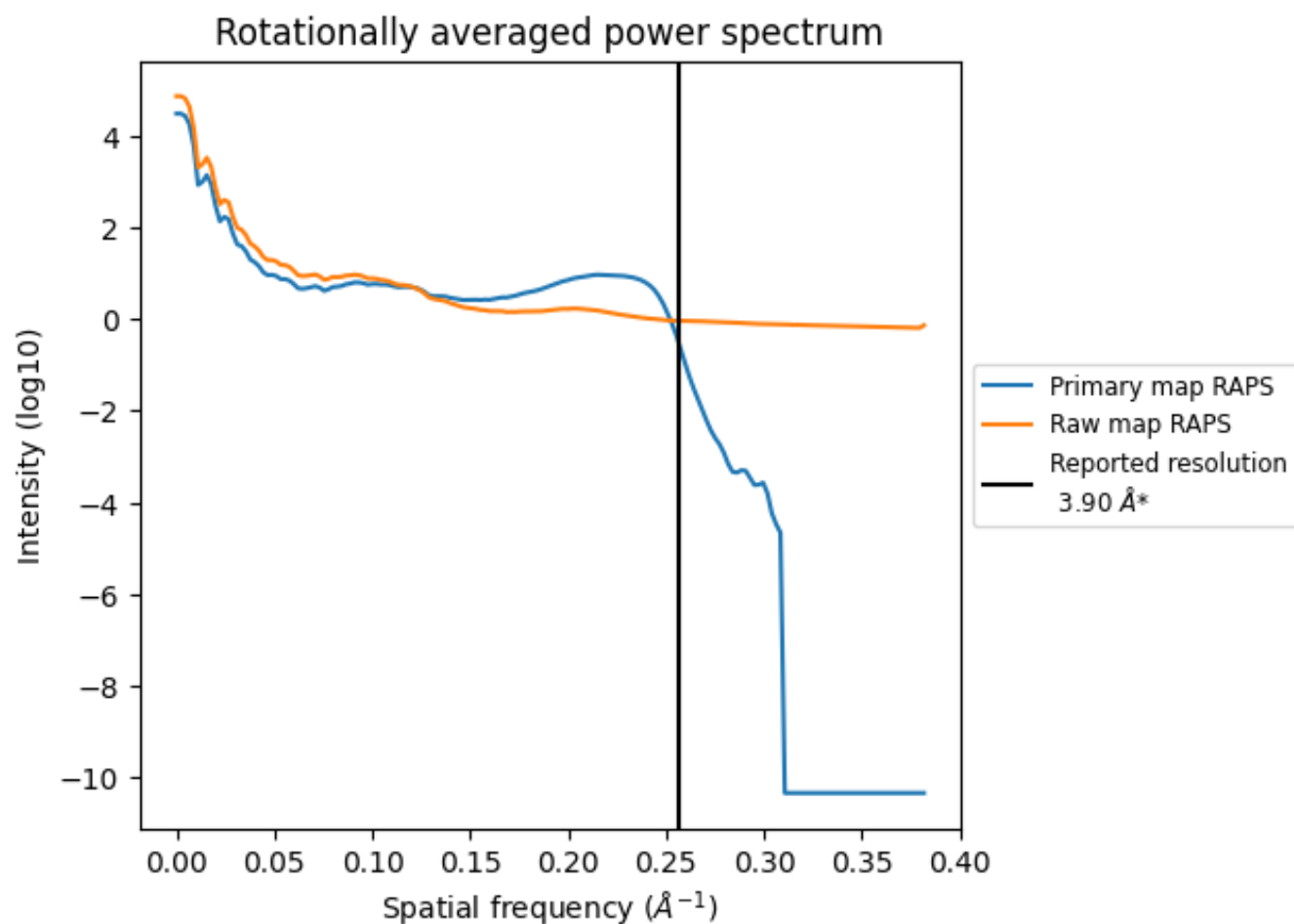
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 307 nm³; this corresponds to an approximate mass of 277 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 ⓘ

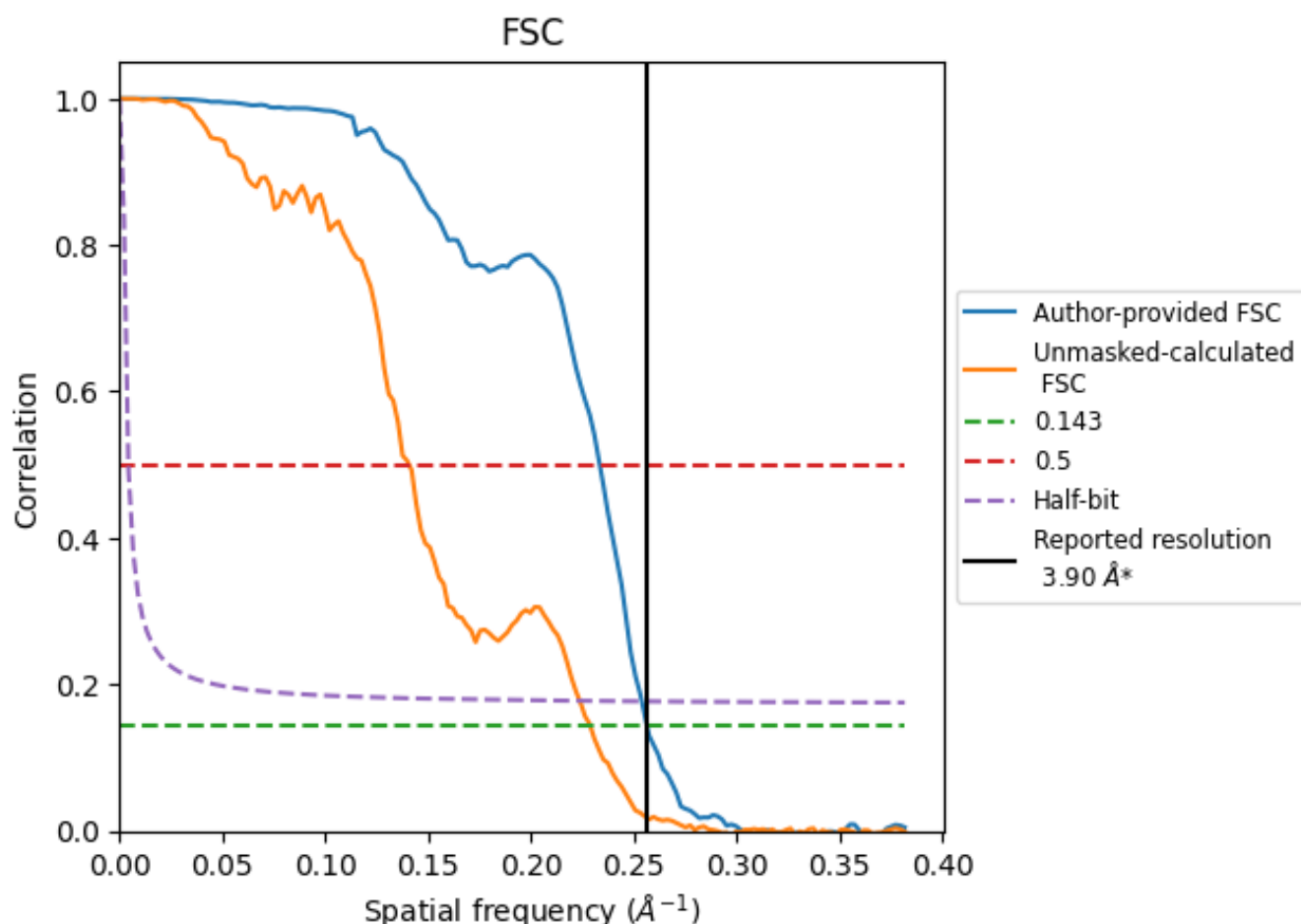


*Reported resolution corresponds to spatial frequency of 0.256 Å⁻¹

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 0.256 Å⁻¹

8.2 Resolution estimates [i](#)

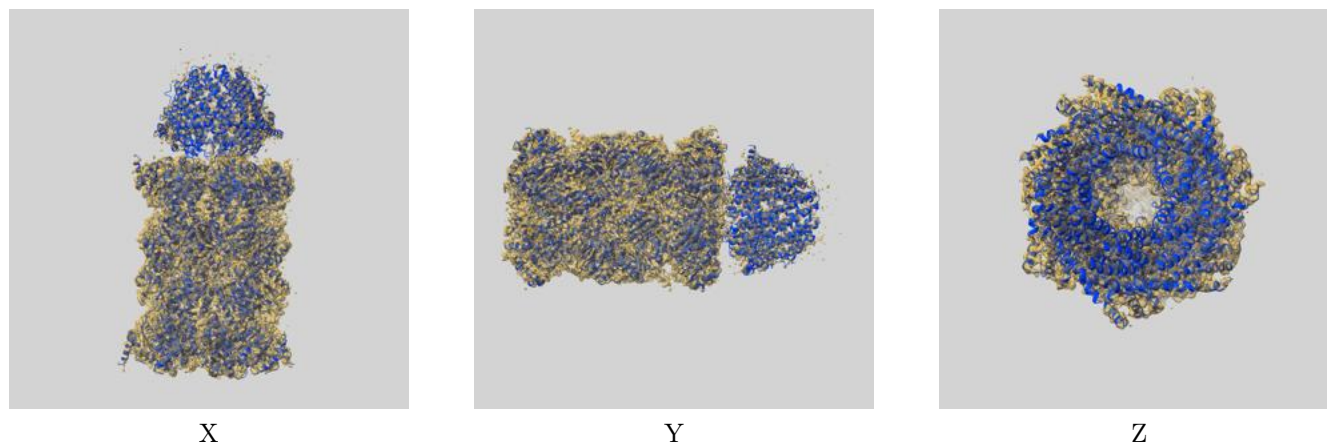
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.90	-	-
Author-provided FSC curve	3.90	4.28	3.94
Unmasked-calculated*	4.37	7.11	4.47

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.37 differs from the reported value 3.9 by more than 10 %

9 Map-model fit [i](#)

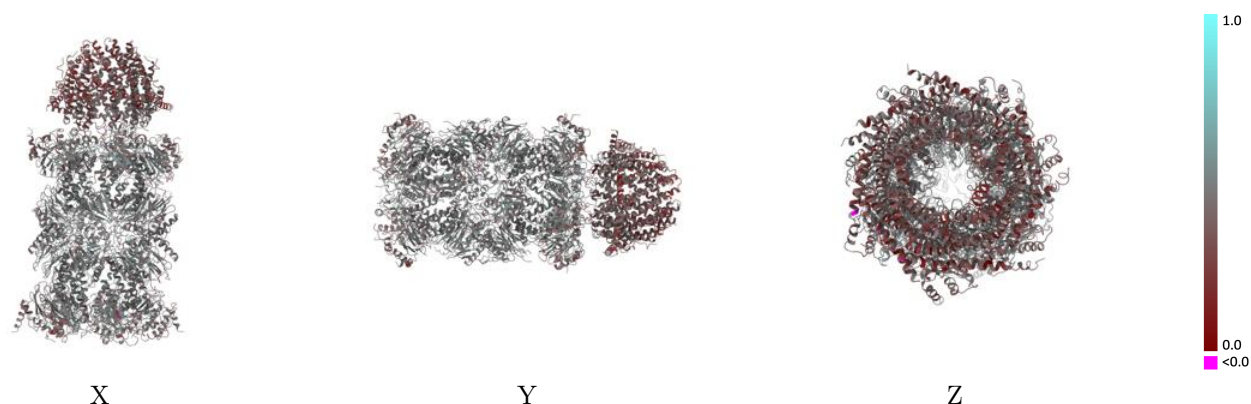
This section contains information regarding the fit between EMDB map EMD-9259 and PDB model 6MUX. Per-residue inclusion information can be found in section [3](#) on page [9](#).

9.1 Map-model overlay [i](#)



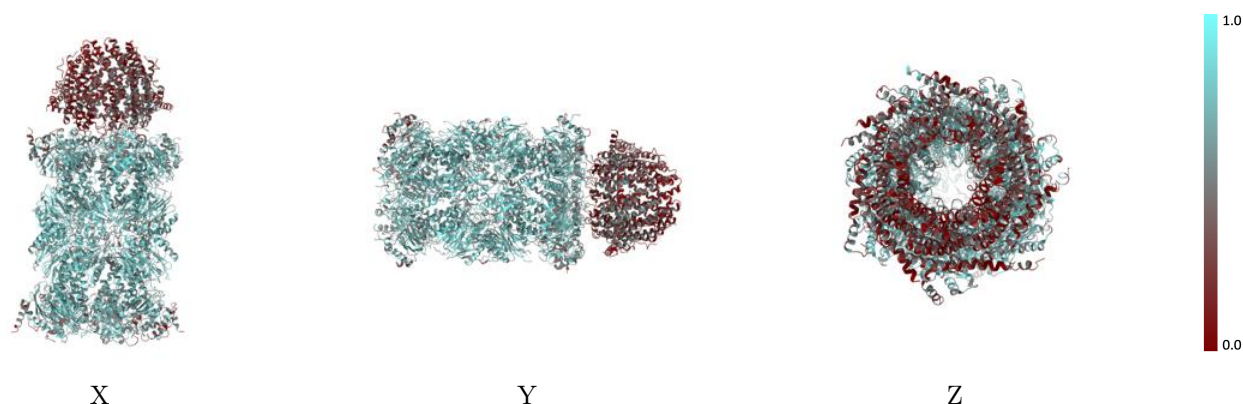
The images above show the 3D surface view of the map at the recommended contour level 6.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



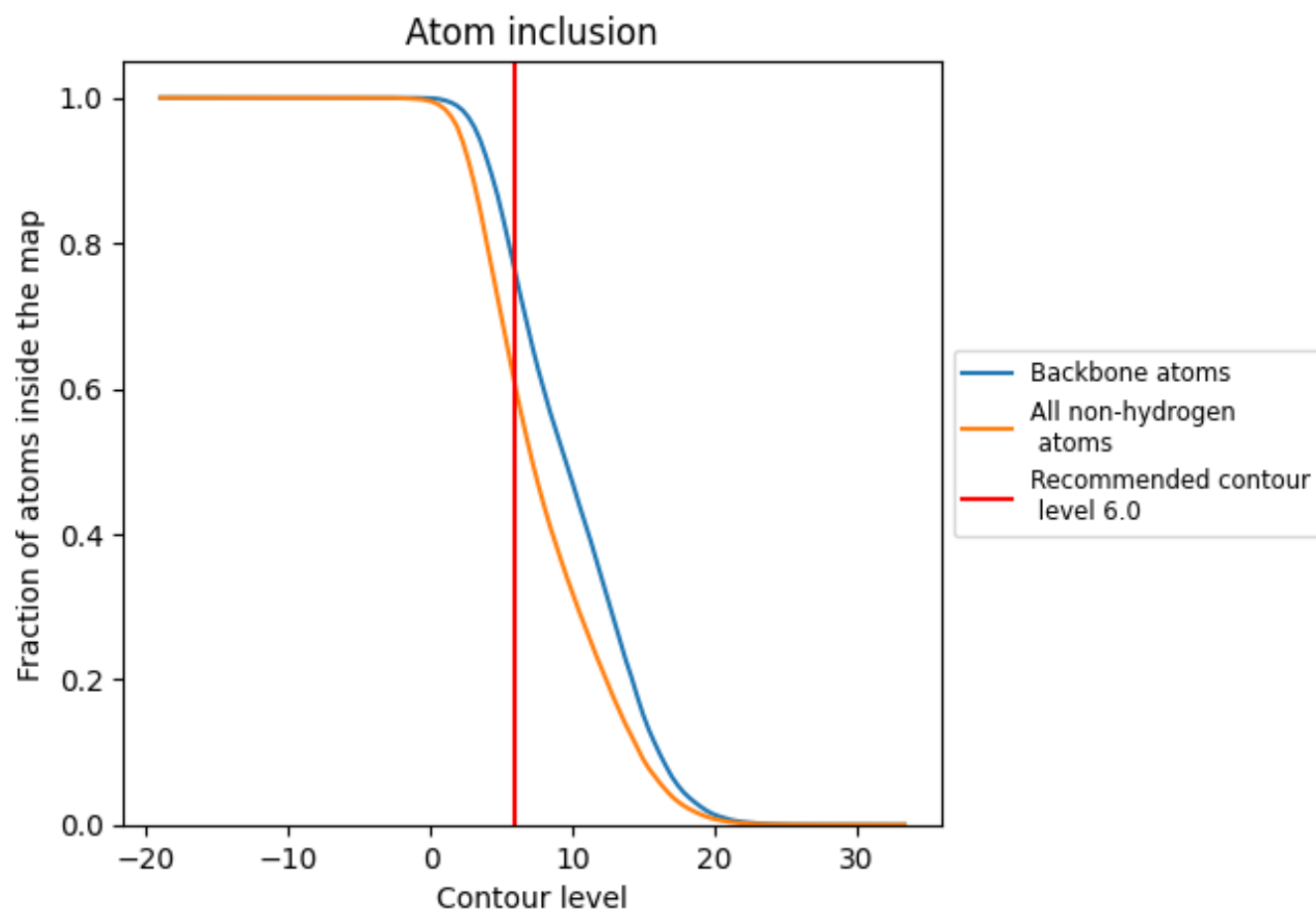
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (6.0).









































































9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 60% 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 (6.0) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6025	 0.4390
A	 0.6436	 0.4620
B	 0.6638	 0.4590
C	 0.6665	 0.4550
D	 0.6237	 0.4520
E	 0.6046	 0.4530
F	 0.6625	 0.4540
G	 0.6961	 0.4640
H	 0.6899	 0.4470
I	 0.6727	 0.4500
J	 0.7516	 0.4800
K	 0.7722	 0.4800
L	 0.7518	 0.4750
M	 0.7213	 0.4800
N	 0.6822	 0.4760
O	 0.5936	 0.4450
P	 0.6340	 0.4550
Q	 0.6691	 0.4520
R	 0.6295	 0.4560
S	 0.6172	 0.4450
T	 0.6605	 0.4500
U	 0.6752	 0.4610
V	 0.6833	 0.4530
W	 0.6709	 0.4560
X	 0.7420	 0.4760
Y	 0.7595	 0.4770
Z	 0.7568	 0.4720
a	 0.7377	 0.4840
b	 0.6994	 0.4760
c	 0.3032	 0.3570
d	 0.2001	 0.3330
e	 0.1734	 0.3040
f	 0.2001	 0.3210
g	 0.2628	 0.3390
h	 0.3146	 0.3550
i	 0.3338	 0.3640

