



wwPDB EM Validation Summary Report ⓘ

Nov 19, 2022 – 09:48 PM EST

PDB ID : 3IZI
EMDB ID : EMD-5245
Title : Mm-cpn rls with ATP
Authors : Douglas, N.R.; Reissmann, S.; Zhang, J.; Chen, B.; Jakana, J.; Kumar, R.;
Chiu, W.; Frydman, J.
Deposited on : 2010-10-29
Resolution : 6.70 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.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) : 2.31.3

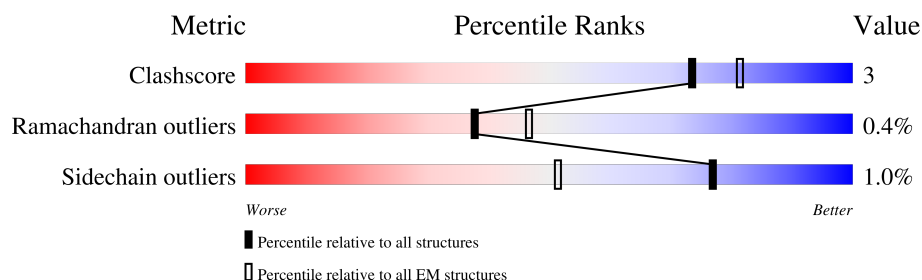
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 6.70 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	513	
1	B	513	
1	C	513	
1	D	513	
1	E	513	
1	F	513	
1	G	513	
1	H	513	

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Mol	Chain	Length	Quality of chain
1	I	513	 38% 90% 9%
1	J	513	 38% 91% 9%
1	K	513	 38% 91% 9%
1	L	513	 38% 91% 9%
1	M	513	 38% 90% 9%
1	N	513	 38% 90% 9%
1	O	513	 38% 89% 10%
1	P	513	 38% 90% 9%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 61440 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 Chaperonin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	B	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	C	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	D	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	E	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	F	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	G	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	H	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	I	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	J	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	K	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	L	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	M	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	N	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	O	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		
1	P	513	Total	C	N	O	S	0	0
			3840	2385	662	768	25		

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	321	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
A	322	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
A	324	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
A	325	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
B	834	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
B	835	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
B	837	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
B	838	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
C	1347	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
C	1348	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
C	1350	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
C	1351	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
D	1860	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
D	1861	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
D	1863	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
D	1864	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
E	2373	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
E	2374	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
E	2376	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
E	2377	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
F	2886	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
F	2887	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
F	2889	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
F	2890	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
G	3399	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
G	3400	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
G	3402	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
G	3403	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
H	3912	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
H	3913	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
H	3915	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
H	3916	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
I	4425	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
I	4426	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
I	4428	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
I	4429	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
J	4938	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
J	4939	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
J	4941	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
J	4942	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
K	5451	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
K	5452	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
K	5454	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8

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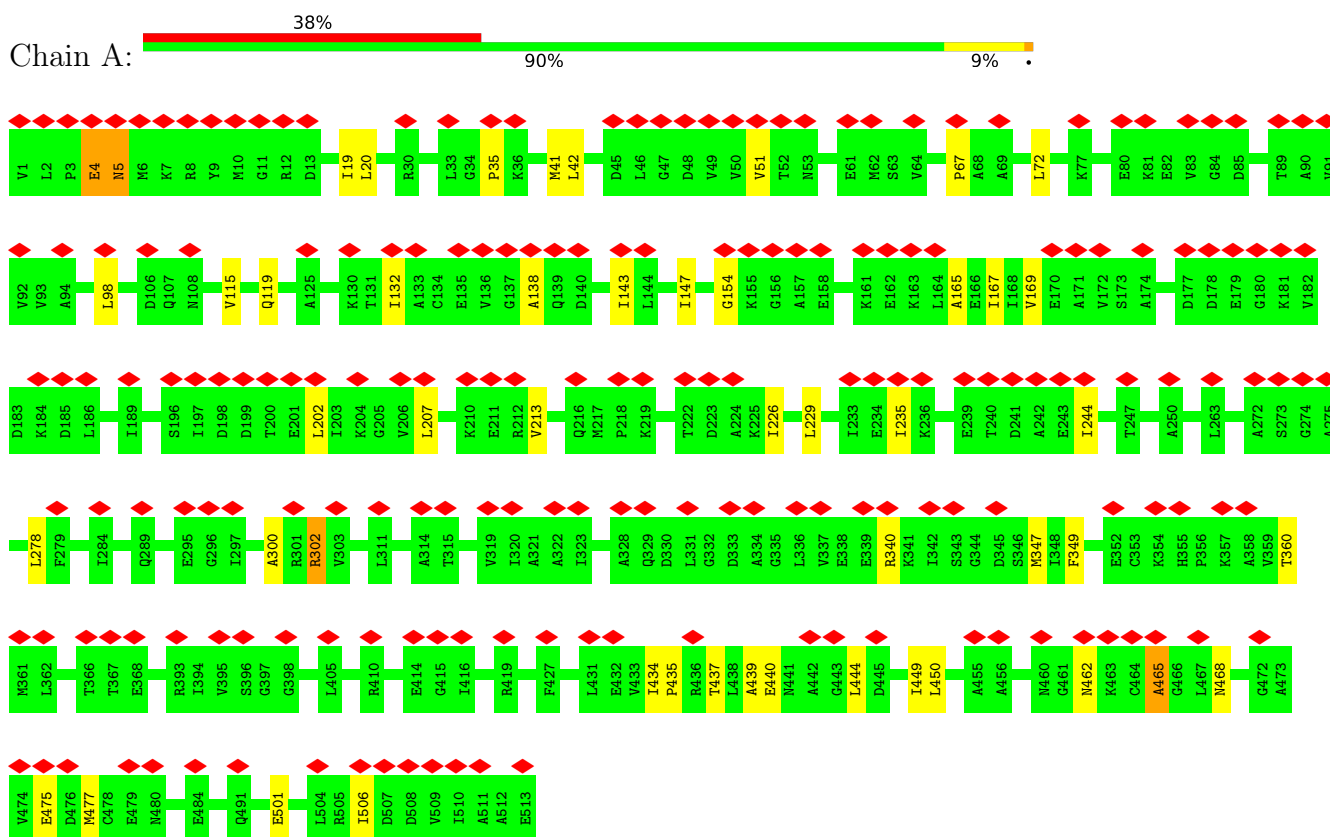
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Chain	Residue	Modelled	Actual	Comment	Reference
K	5455	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
L	5964	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
L	5965	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
L	5967	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
L	5968	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
M	6477	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
M	6478	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
M	6480	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
M	6481	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
N	6990	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
N	6991	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
N	6993	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
N	6994	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
O	7503	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
O	7504	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
O	7506	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
O	7507	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8
P	8016	ALA	THR	ENGINEERED MUTATION	UNP Q877G8
P	8017	ALA	ASN	ENGINEERED MUTATION	UNP Q877G8
P	8019	ALA	LYS	ENGINEERED MUTATION	UNP Q877G8
P	8020	ALA	ASP	ENGINEERED MUTATION	UNP Q877G8

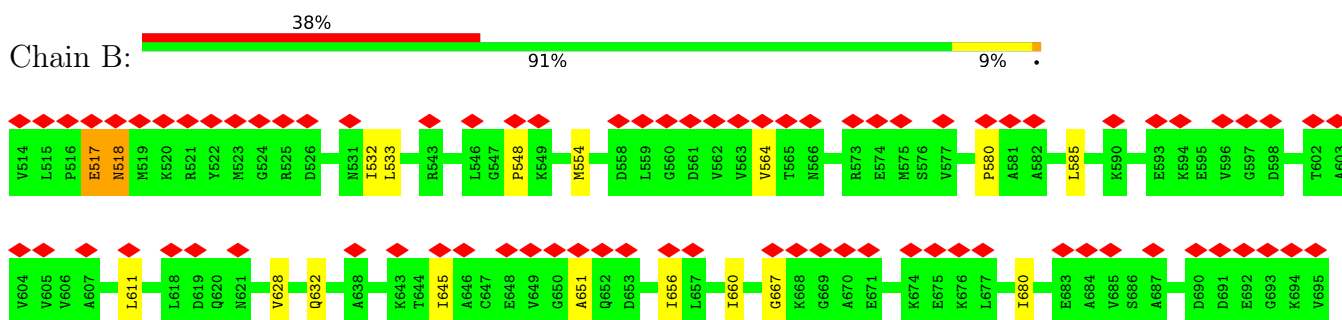
3 Residue-property plots

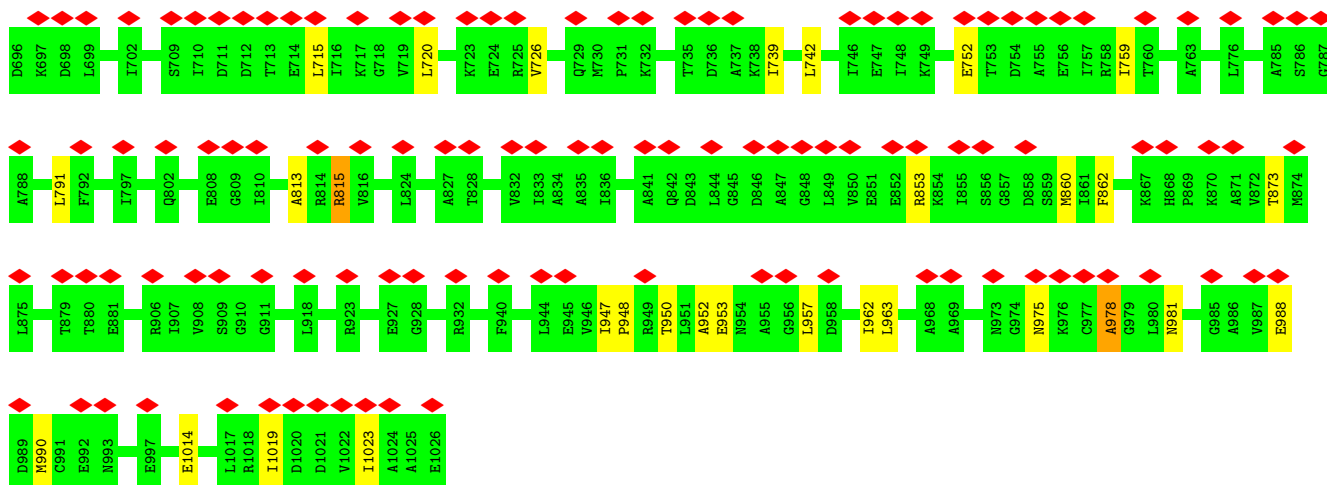
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Chaperonin

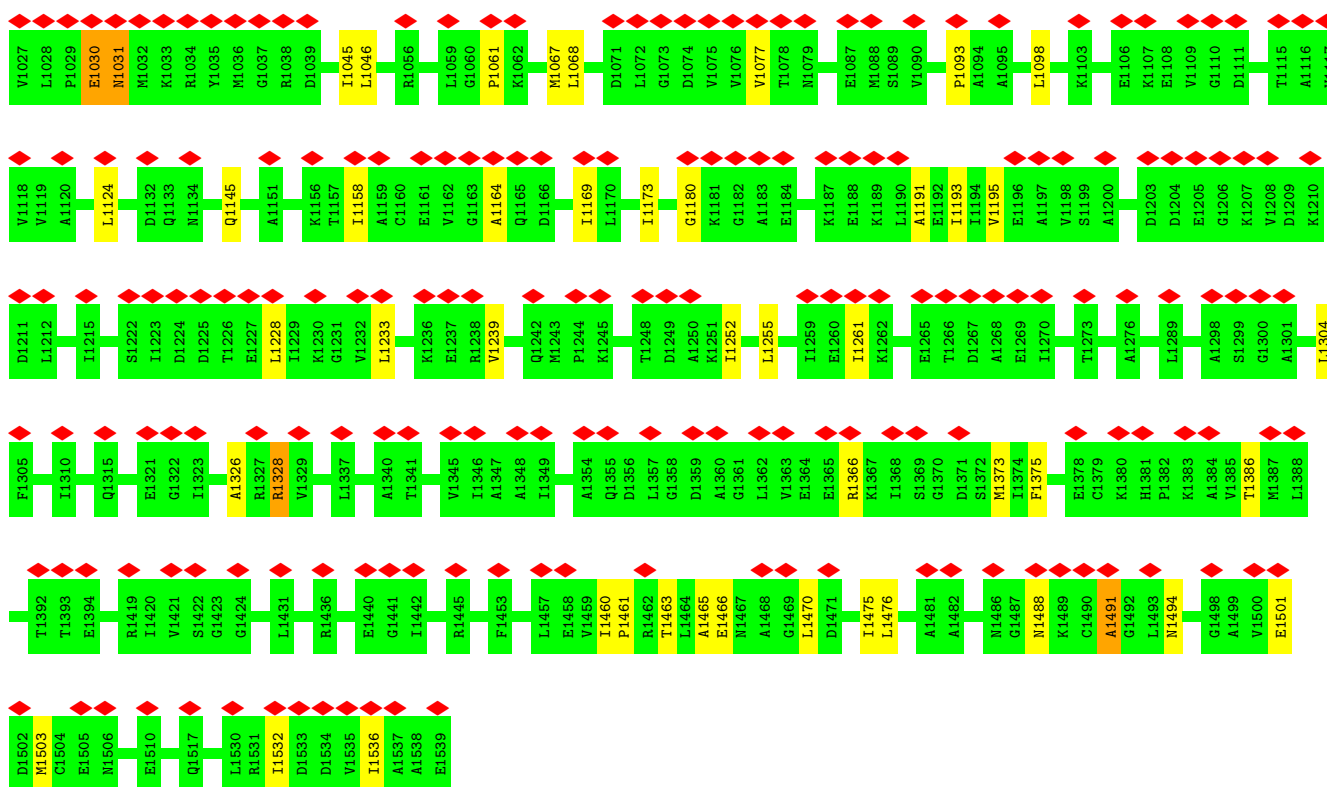
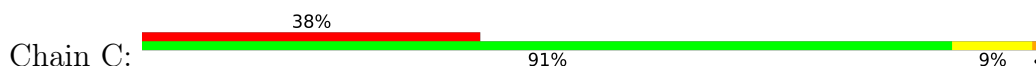


• Molecule 1: Chaperonin

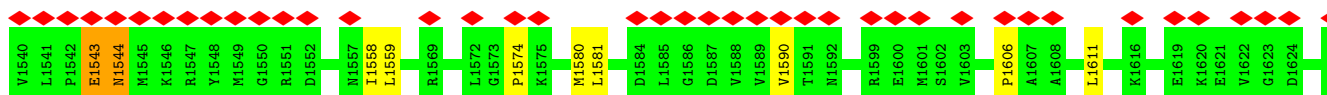
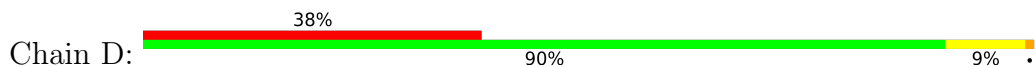


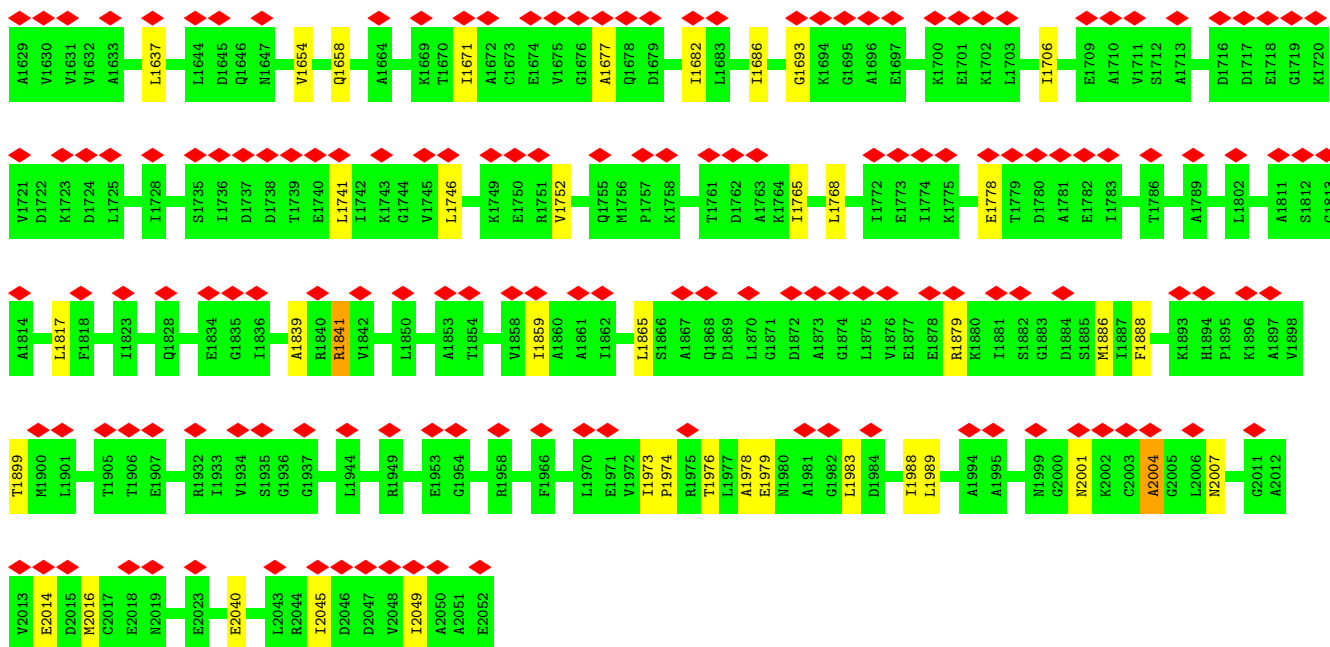


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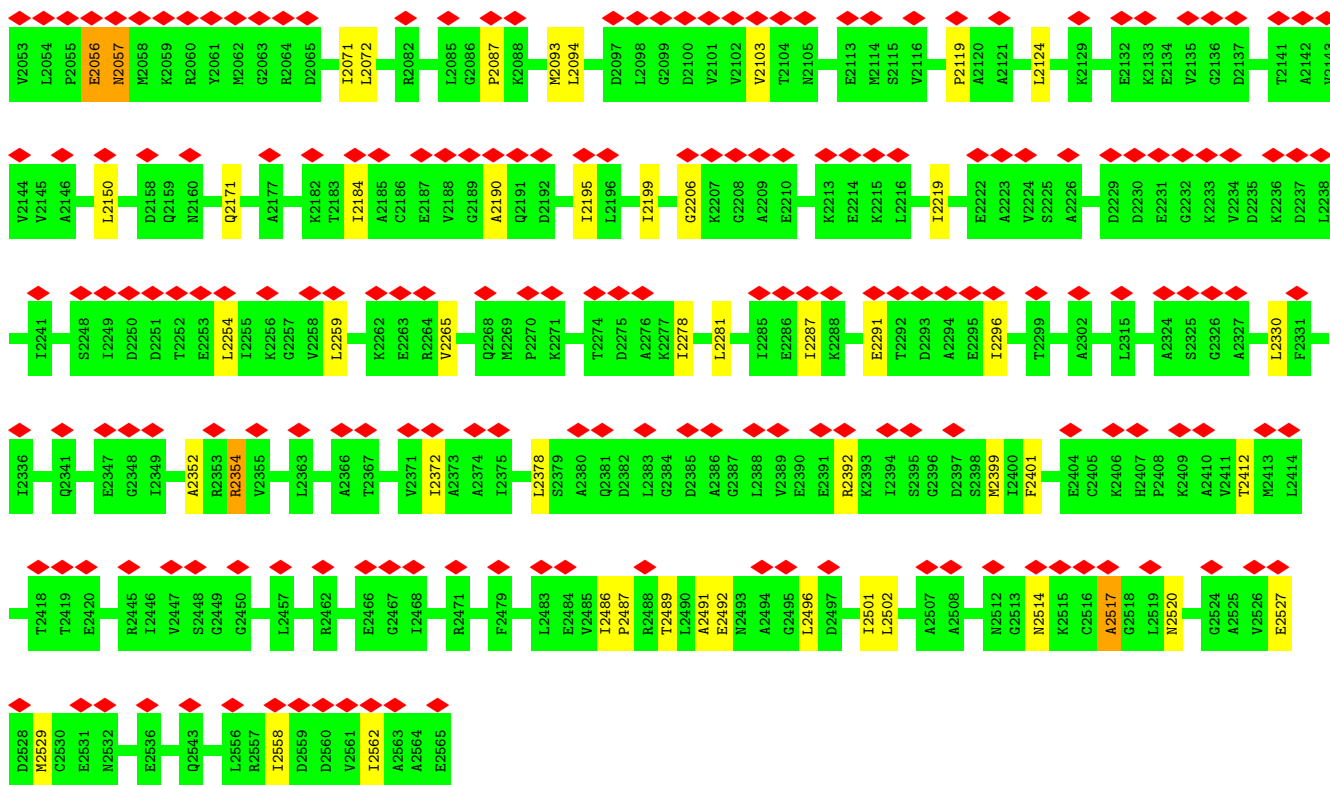
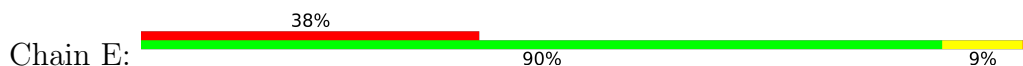


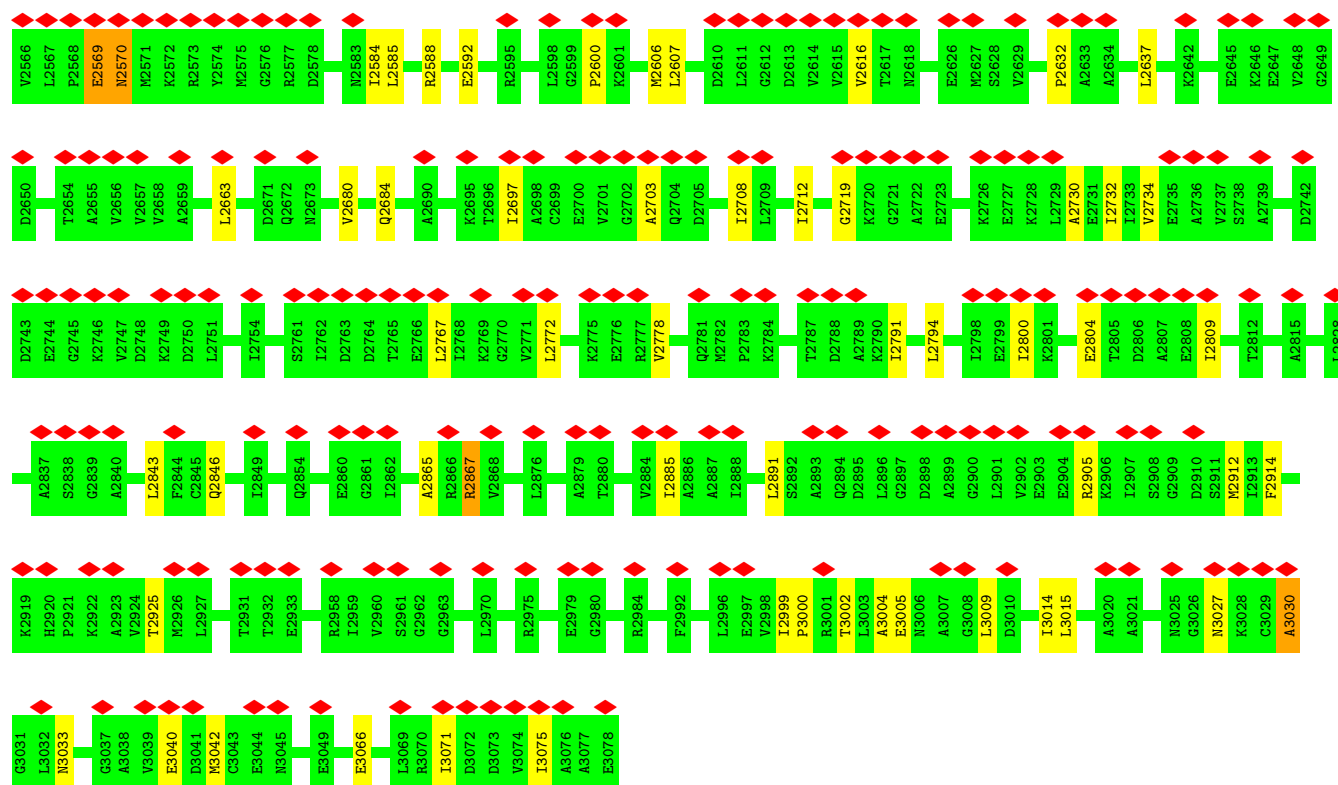
• Molecule 1: Chaperonin



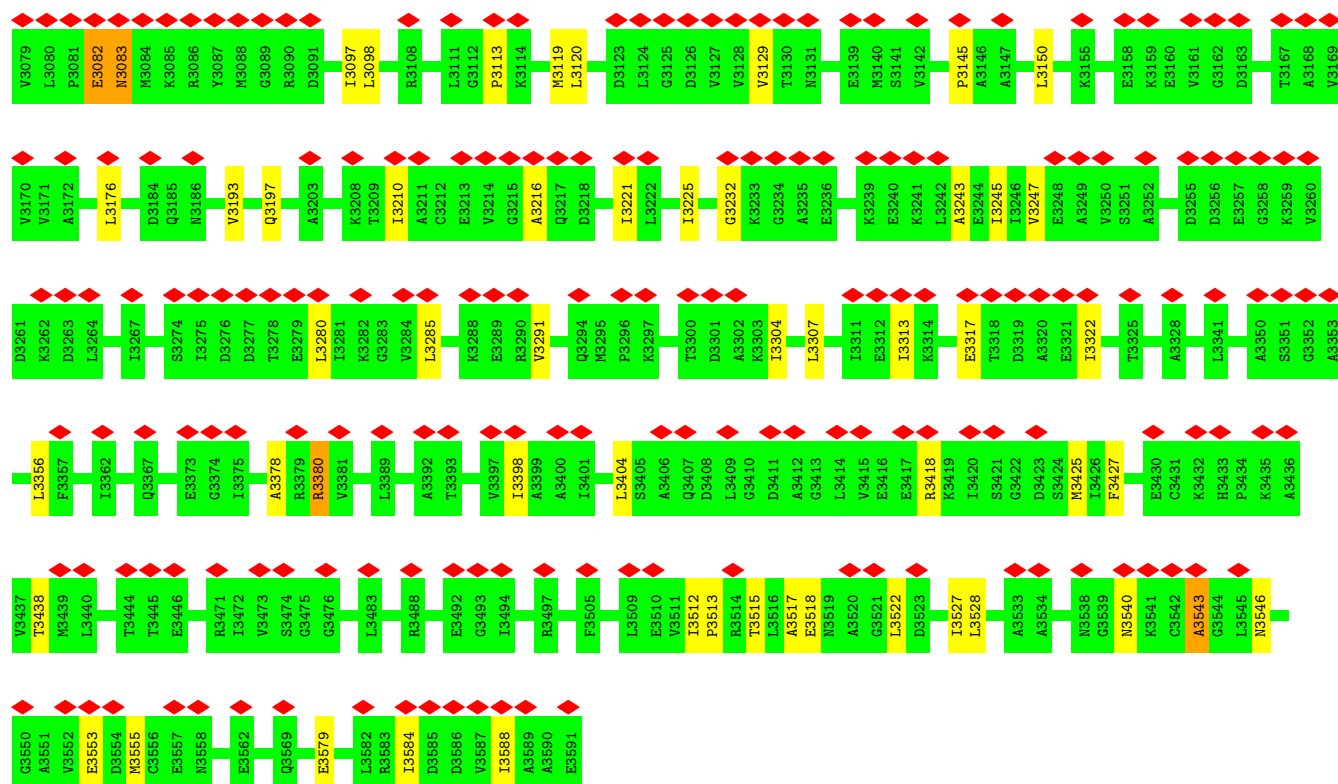
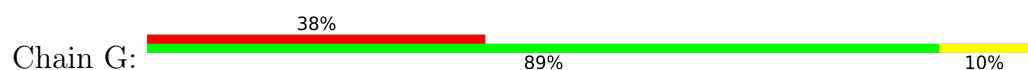


• Molecule 1: Chaperonin



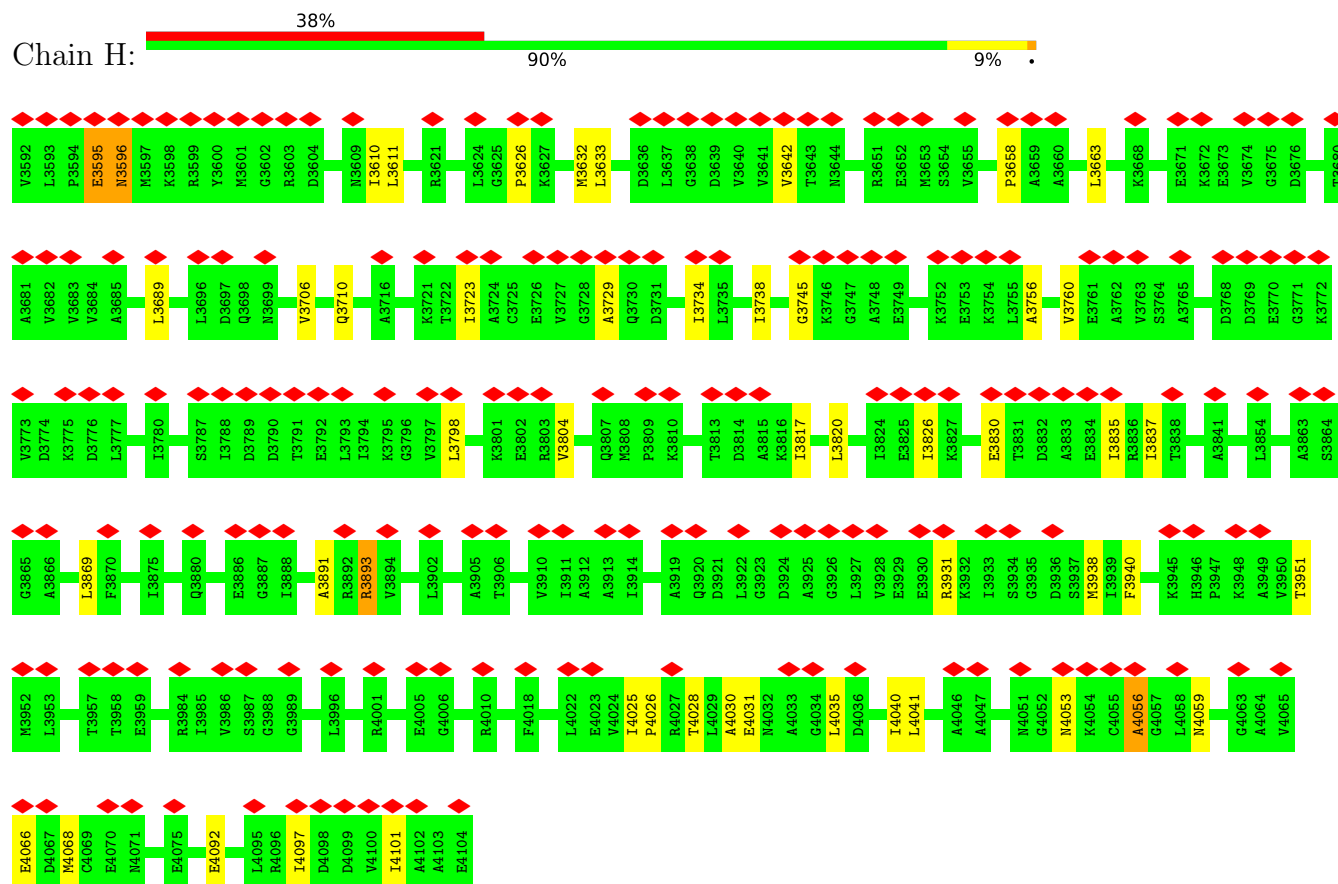


• Molecule 1: Chaperonin



• Molecule 1: Chaperonin

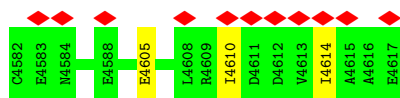
Chain H:



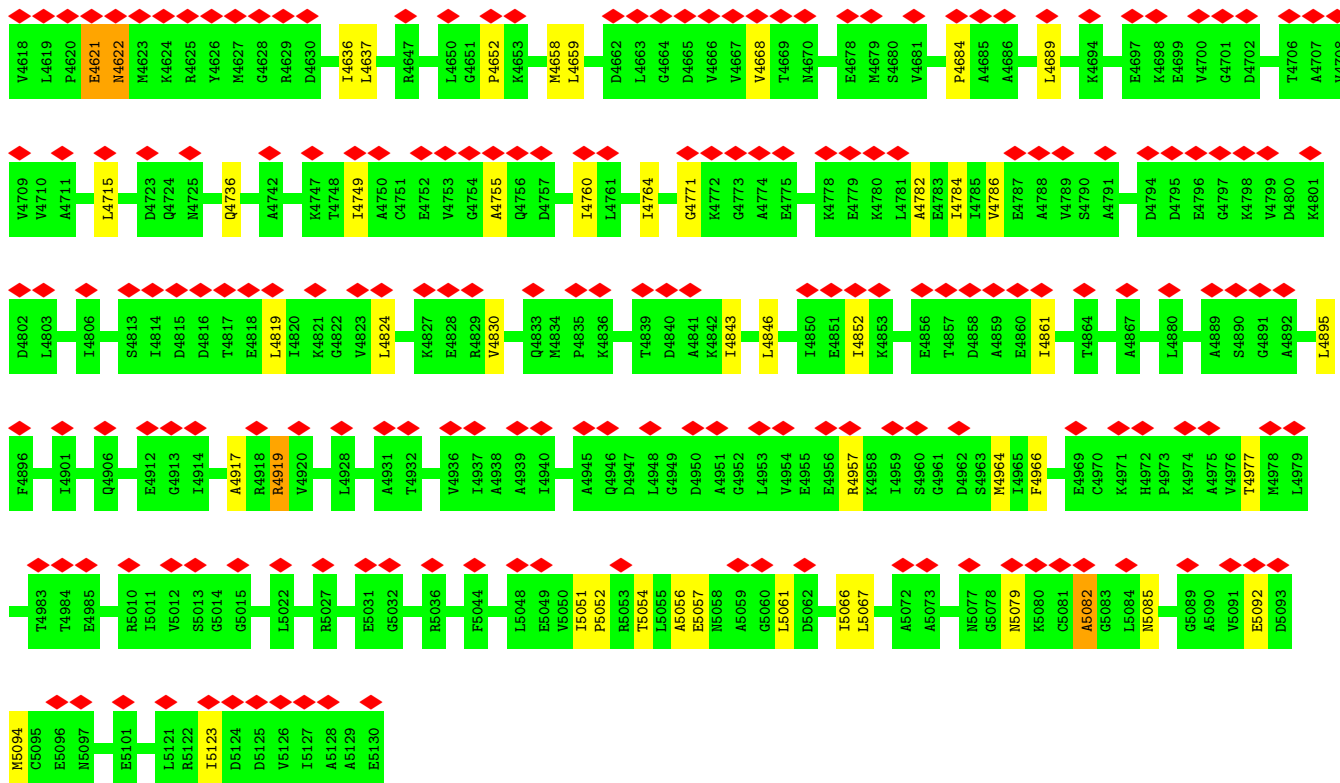
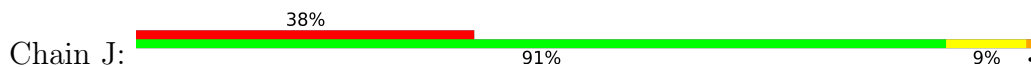
• Molecule 1: Chaperonin

Chain I:

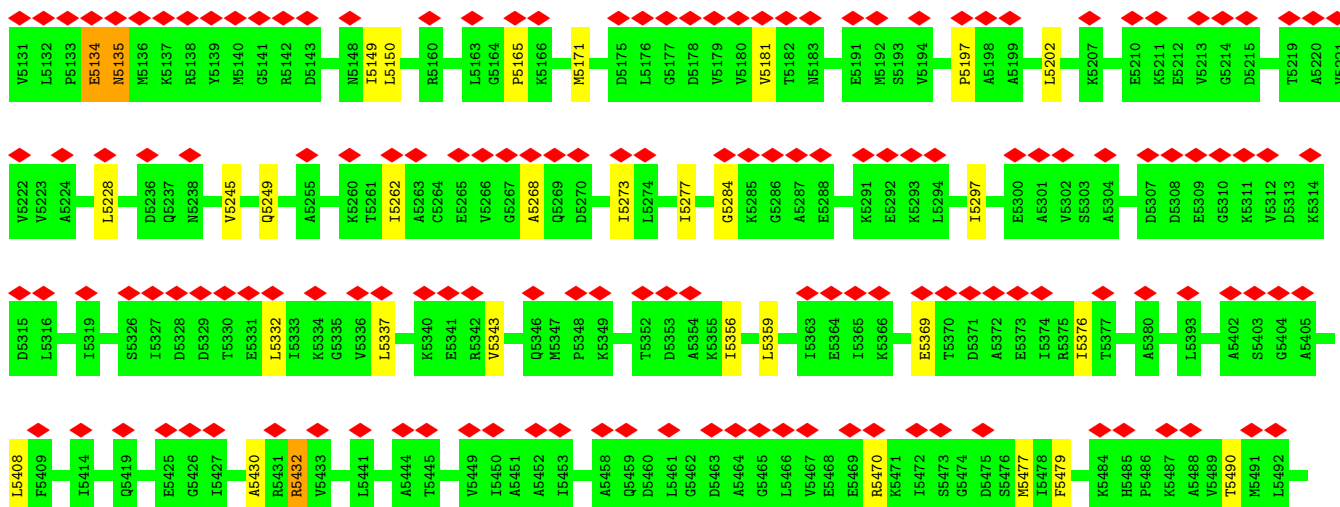
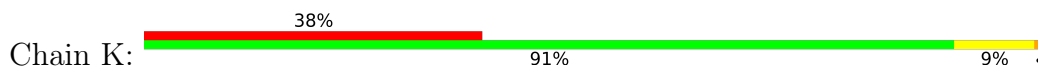


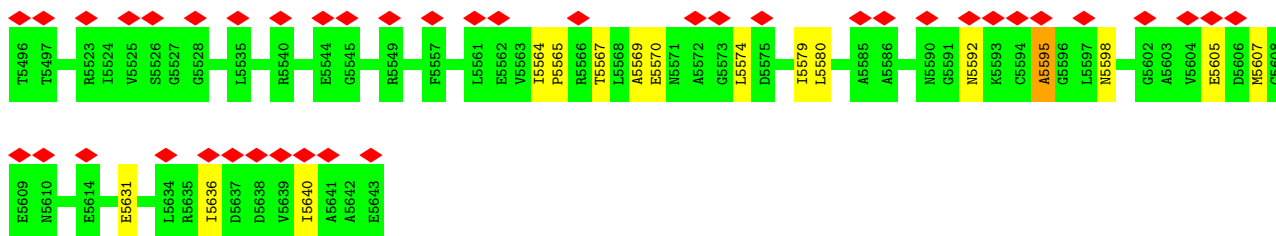


• Molecule 1: Chaperonin

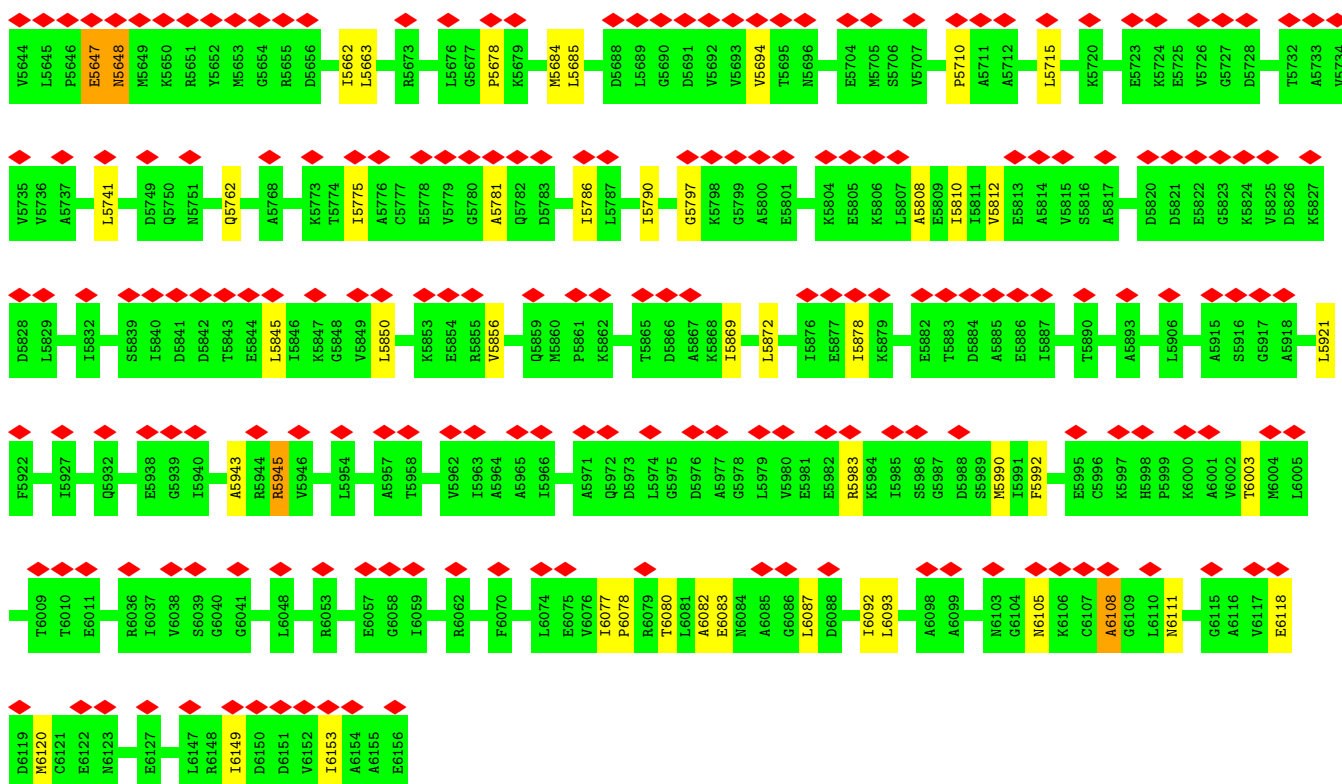
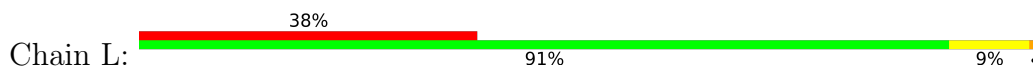


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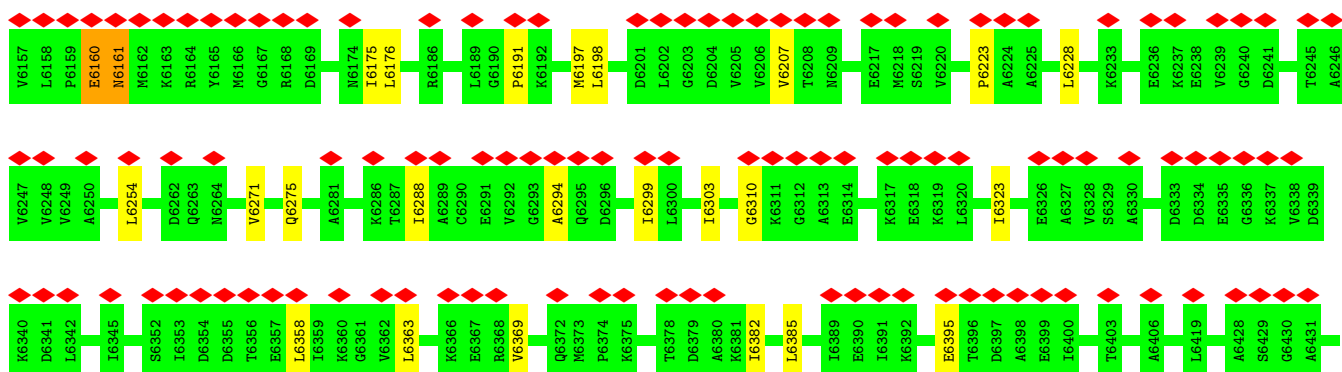
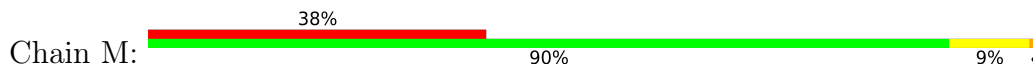


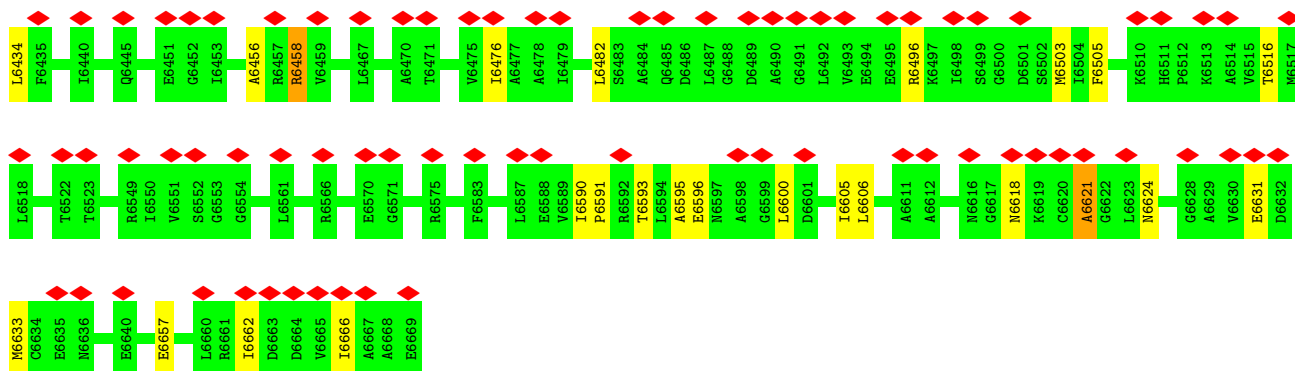


• Molecule 1: Chaperonin

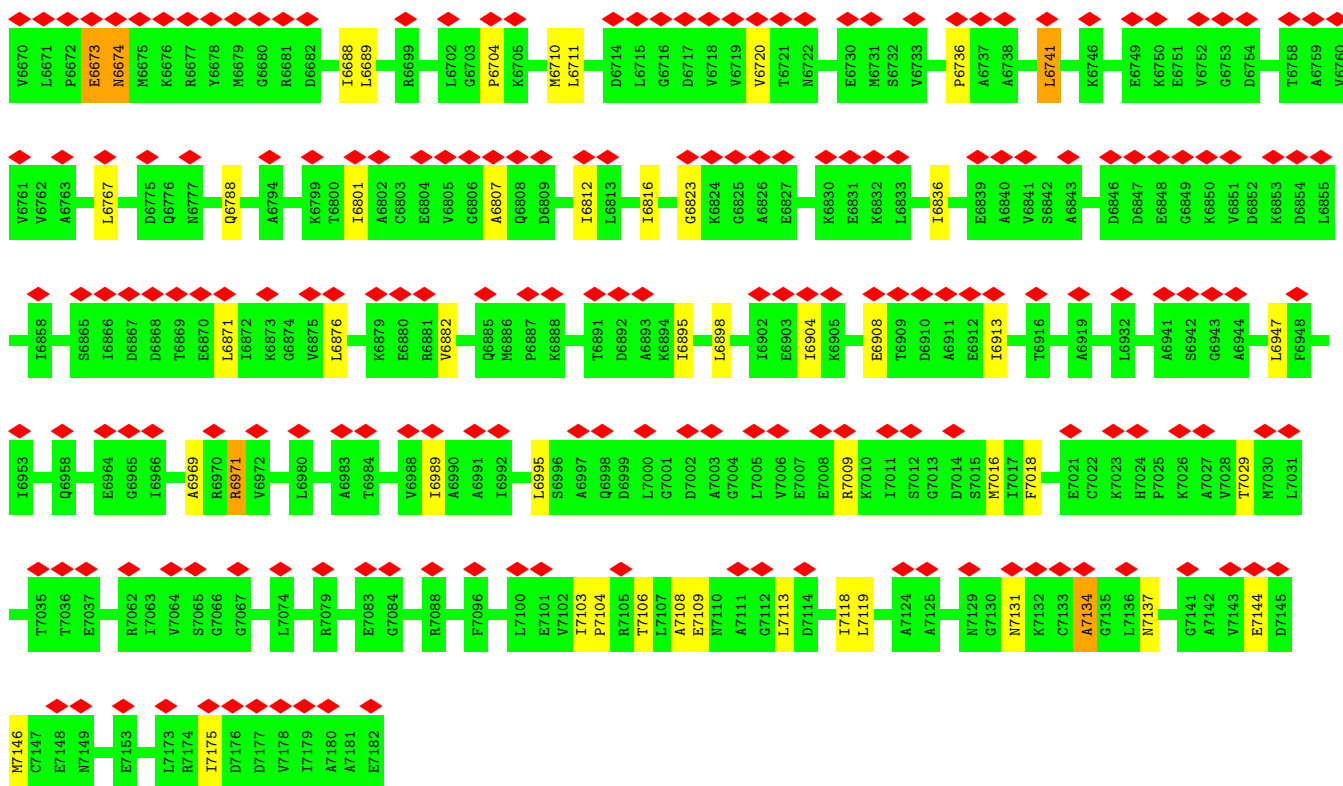
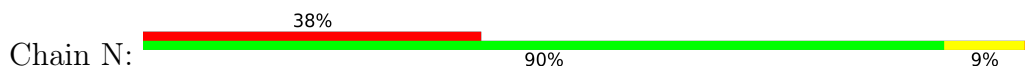


• Molecule 1: Chaperonin

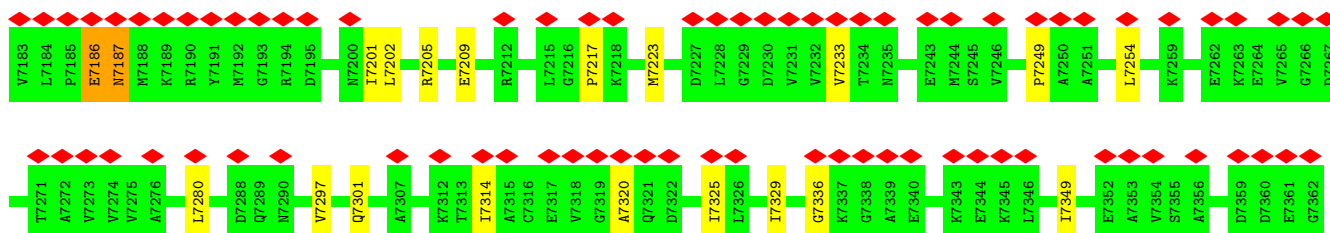
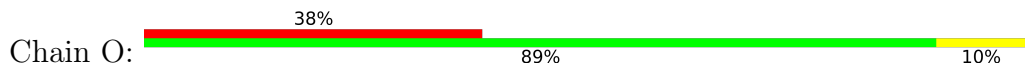


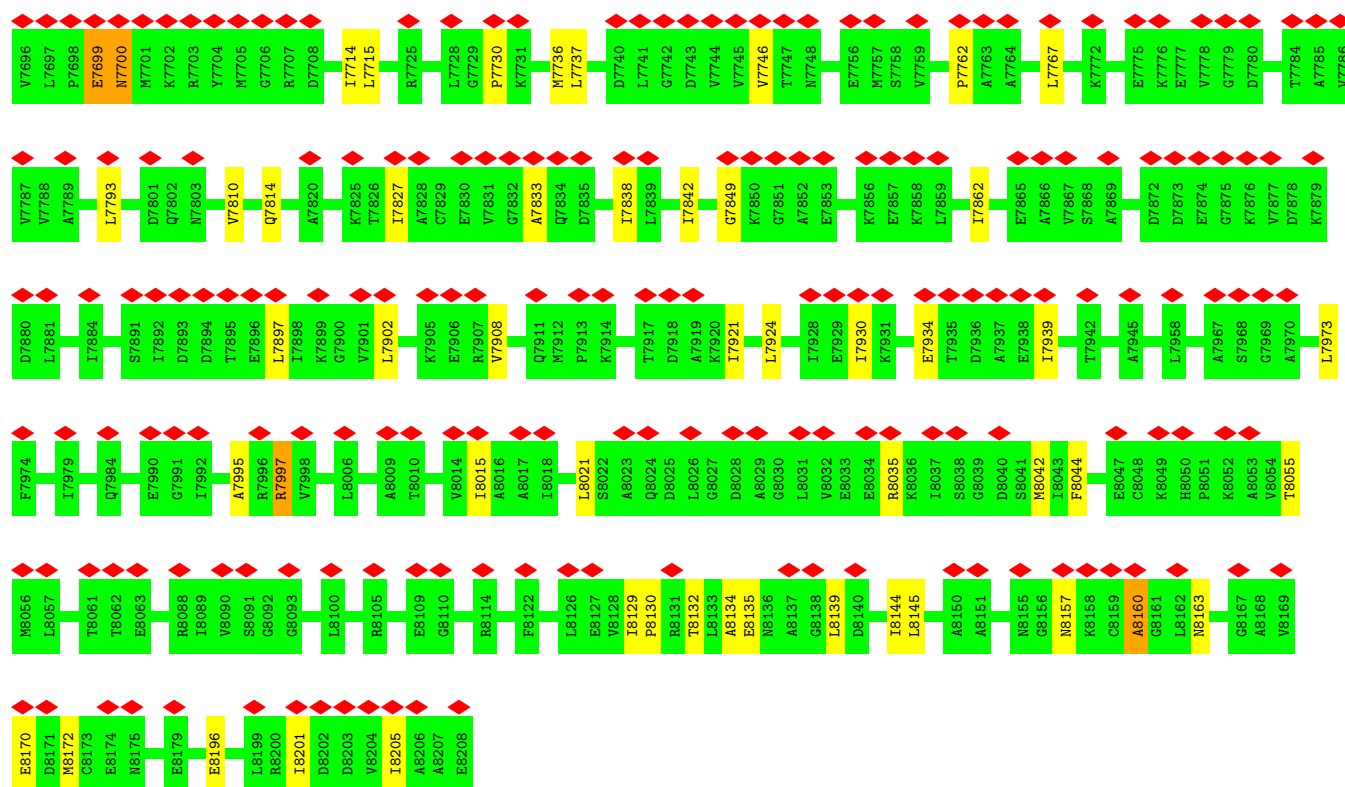


• Molecule 1: Chaperonin



• Molecule 1: Chaperonin





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, D8	Depositor
Number of particles used	Not provided	
Resolution determination method	FSC 0.5 CUT-OFF	Depositor
CTF correction method	Not provided	
Microscope	JEOL 3200FSC	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN ULTRASCAN 4000 (4k x 4k)	Depositor
Maximum map value	0.744	Depositor
Minimum map value	-0.366	Depositor
Average map value	0.017	Depositor
Map value standard deviation	0.071	Depositor
Recommended contour level	0.28	Depositor
Map size (\AA)	255.36002, 255.36002, 255.36002	wwPDB
Map dimensions	192, 192, 192	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.33, 1.33, 1.33	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	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	B	0.88	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	C	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	D	0.88	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	E	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	F	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	G	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	H	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	I	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	J	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	K	0.88	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	L	0.88	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	M	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	N	0.88	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	O	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
1	P	0.87	2/3863 (0.1%)	0.78	2/5199 (0.0%)
All	All	0.87	32/61808 (0.1%)	0.78	32/83184 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	3
1	C	0	3
1	D	0	3
1	E	0	3
1	F	0	3
1	G	0	3
1	H	0	3
1	I	0	3

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Mol	Chain	#Chirality outliers	#Planarity outliers
1	J	0	3
1	K	0	3
1	L	0	3
1	M	0	3
1	N	0	3
1	O	0	3
1	P	0	3
All	All	0	48

The worst 5 of 32 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	2056	GLU	C-O	-6.03	1.11	1.23
1	D	1543	GLU	C-O	-5.98	1.11	1.23
1	P	7699	GLU	C-O	-5.97	1.12	1.23
1	J	4621	GLU	C-O	-5.97	1.12	1.23
1	O	7186	GLU	C-O	-5.96	1.12	1.23

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	5872	LEU	CB-CG-CD2	-5.29	102.02	111.00
1	O	7411	LEU	CB-CG-CD2	-5.28	102.03	111.00
1	A	229	LEU	CB-CG-CD2	-5.27	102.03	111.00
1	F	2794	LEU	CB-CG-CD2	-5.27	102.04	111.00
1	H	3820	LEU	CB-CG-CD2	-5.27	102.04	111.00

There are no chirality outliers.

5 of 48 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	119	GLN	Mainchain
1	A	138	ALA	Mainchain
1	A	475	GLU	Mainchain
1	B	632	GLN	Mainchain
1	B	651	ALA	Mainchain

5.2 Too-close contacts [i](#)

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	3840	0	3986	30	0
1	B	3840	0	3983	28	0
1	C	3840	0	3983	28	0
1	D	3840	0	3983	30	0
1	E	3840	0	3983	31	0
1	F	3840	0	3983	33	0
1	G	3840	0	3983	32	0
1	H	3840	0	3983	31	0
1	I	3840	0	3983	31	0
1	J	3840	0	3983	29	0
1	K	3840	0	3983	28	0
1	L	3840	0	3983	28	0
1	M	3840	0	3983	29	0
1	N	3840	0	3983	31	0
1	O	3840	0	3983	32	0
1	P	3840	0	3983	31	0
All	All	61440	0	63731	407	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 3.

The worst 5 of 407 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)
1:G:3380:ARG:HE	1:G:3380:ARG:HA	1.60	0.67
1:A:302:ARG:HE	1:A:302:ARG:HA	1.60	0.67
1:P:7997:ARG:HE	1:P:7997:ARG:HA	1.60	0.67
1:J:4919:ARG:HE	1:J:4919:ARG:HA	1.60	0.67
1:H:3893:ARG:HE	1:H:3893:ARG:HA	1.60	0.66

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	B	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	C	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	D	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	E	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	F	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	G	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	H	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	I	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	J	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	K	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	L	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	M	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	N	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	O	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
1	P	511/513 (100%)	493 (96%)	16 (3%)	2 (0%)	34	72
All	All	8176/8208 (100%)	7888 (96%)	256 (3%)	32 (0%)	38	72

5 of 32 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	B	518	ASN
1	C	1031	ASN
1	D	1544	ASN
1	E	2057	ASN

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	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	B	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	C	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	D	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	E	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	F	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	G	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	H	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	I	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	J	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	K	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	L	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	M	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	N	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	O	410/410 (100%)	406 (99%)	4 (1%)	76	86
1	P	410/410 (100%)	406 (99%)	4 (1%)	76	86
All	All	6560/6560 (100%)	6496 (99%)	64 (1%)	77	86

5 of 64 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	O	7254	LEU
1	O	7529	MET
1	G	3098	LEU
1	F	2912	MET
1	P	7715	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
1	O	7289	GLN
1	O	7650	ASN
1	G	3185	GLN
1	F	3033	ASN
1	P	7802	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.

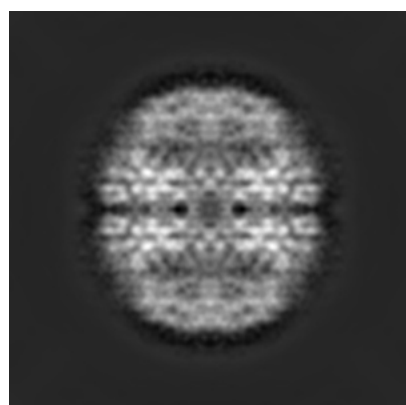
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-5245. These allow visual inspection of the internal detail of the map and identification of artifacts.

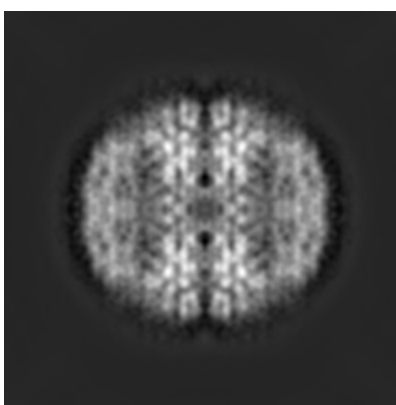
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

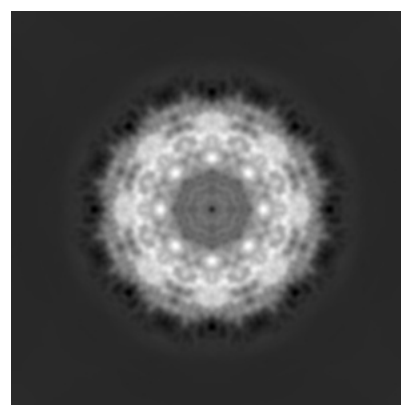
6.1.1 Primary map



X



Y

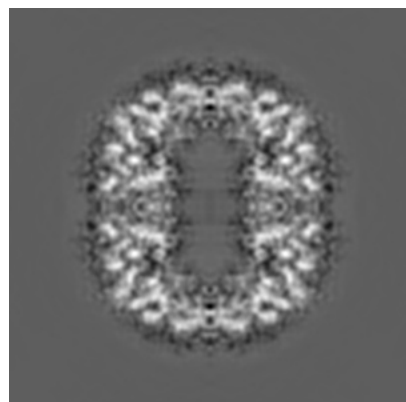


Z

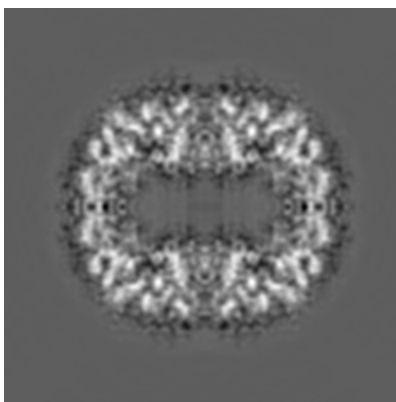
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

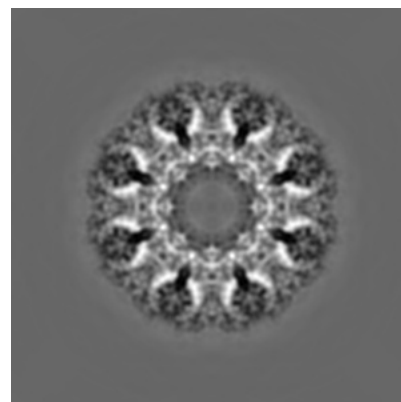
6.2.1 Primary map



X Index: 96



Y Index: 96

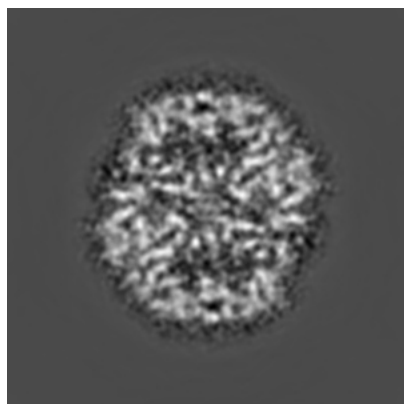


Z Index: 96

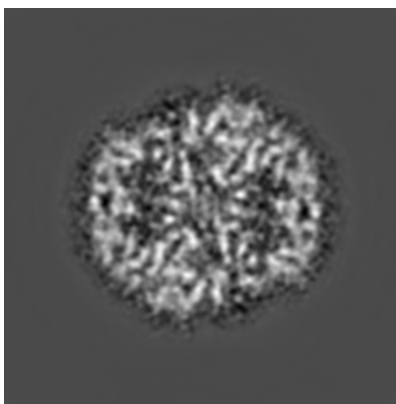
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 69



Y Index: 123

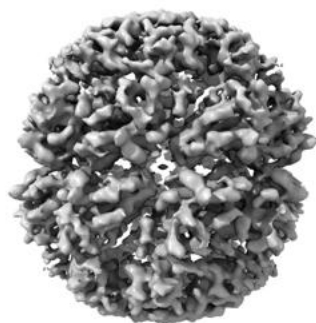


Z Index: 106

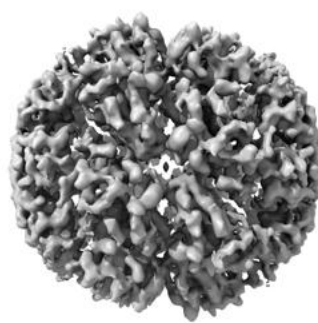
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views [i](#)

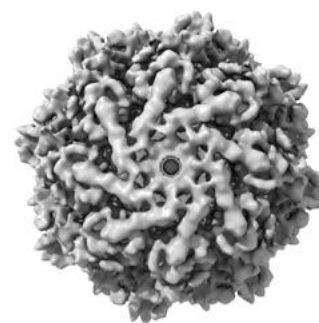
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.28. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

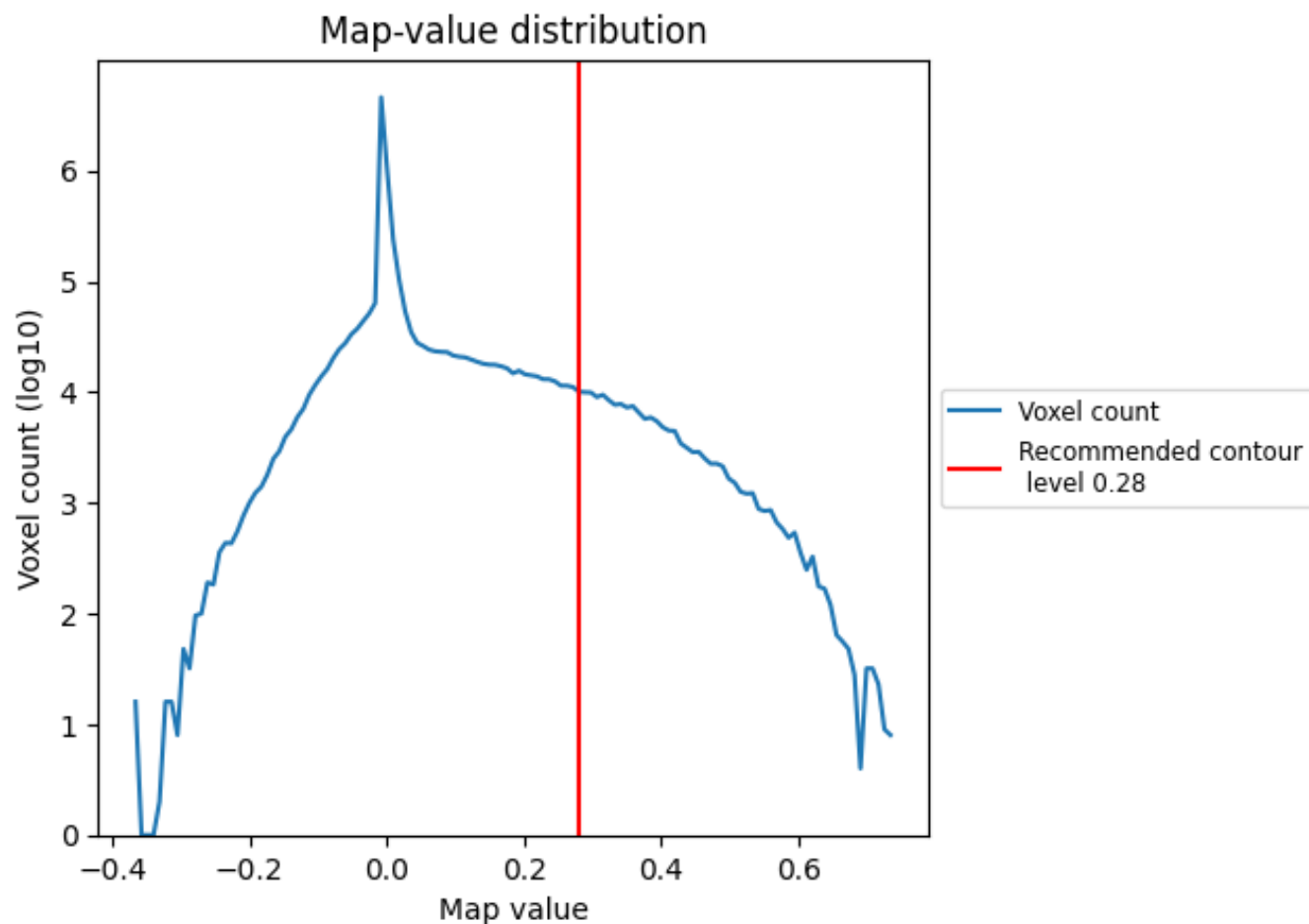
6.5 Mask visualisation

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

7 Map analysis [i](#)

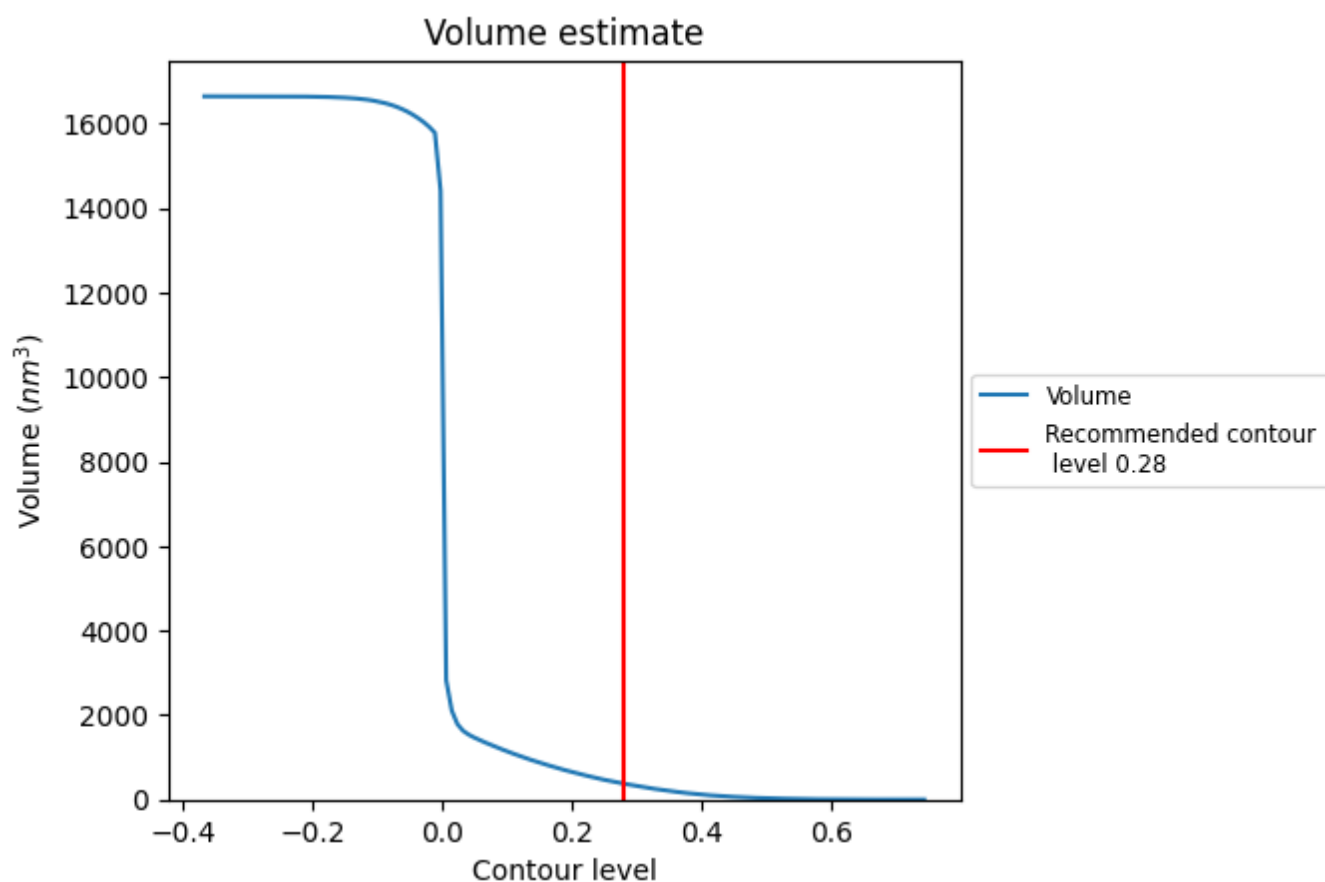
This section contains the results of statistical analysis of the map.

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.

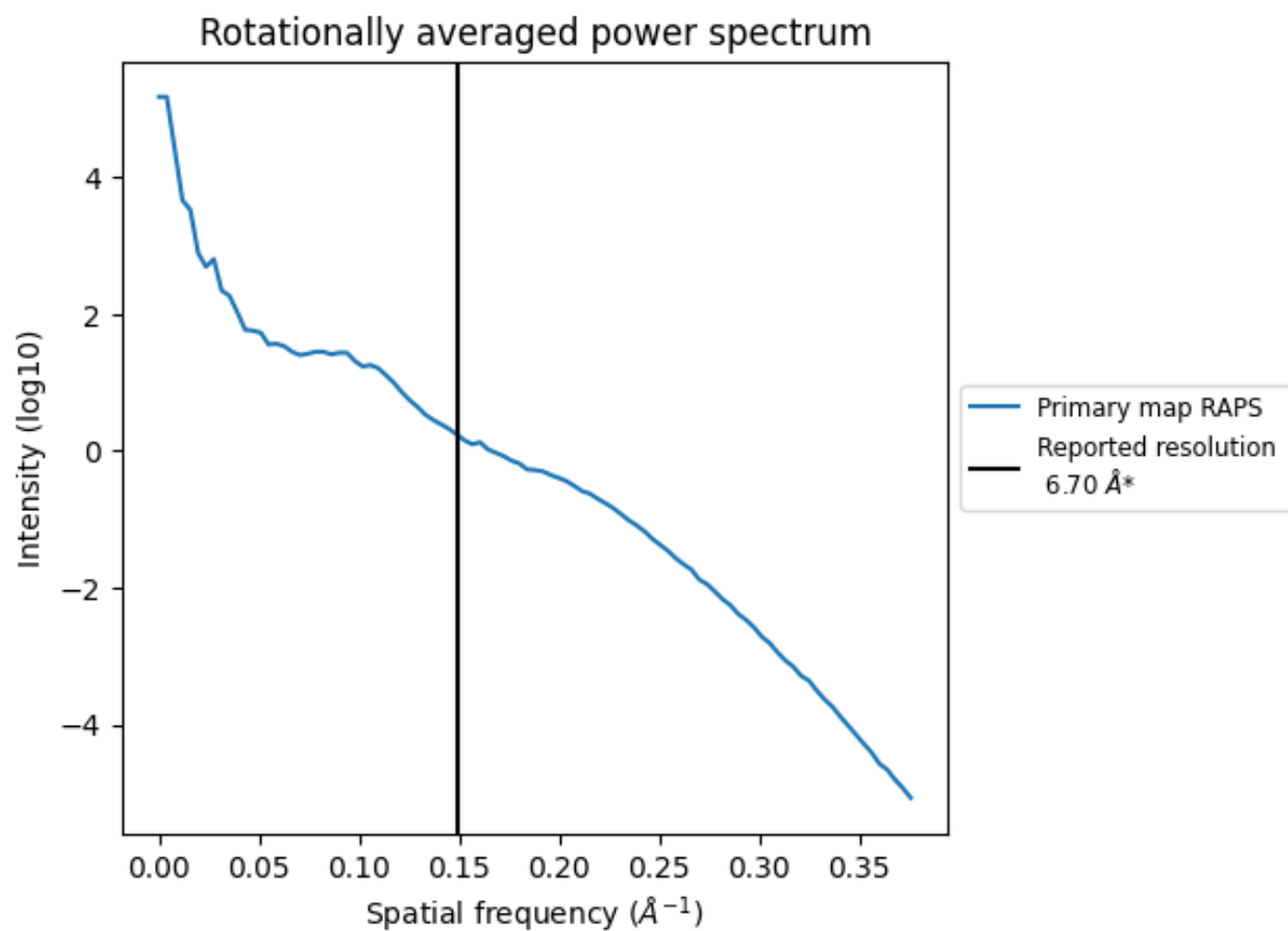
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 379 nm³; this corresponds to an approximate mass of 342 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 [i](#)



*Reported resolution corresponds to spatial frequency of 0.149 Å⁻¹

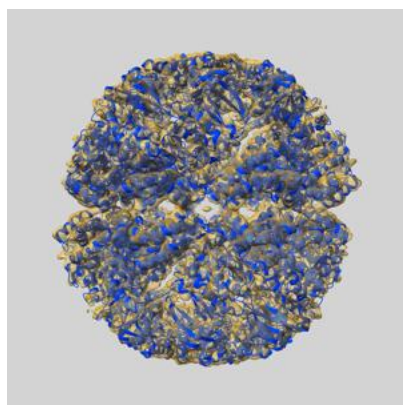
8 Fourier-Shell correlation

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

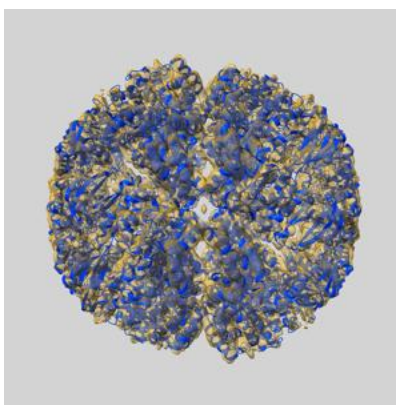
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-5245 and PDB model 3IZI. Per-residue inclusion information can be found in section [3](#) on page [7](#).

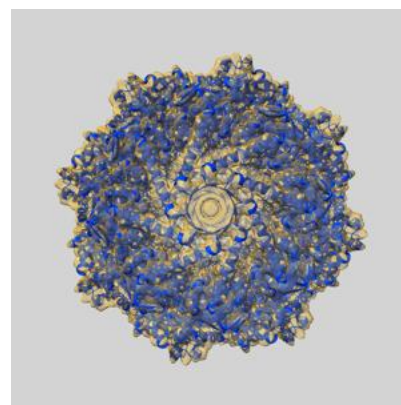
9.1 Map-model overlay [i](#)



X



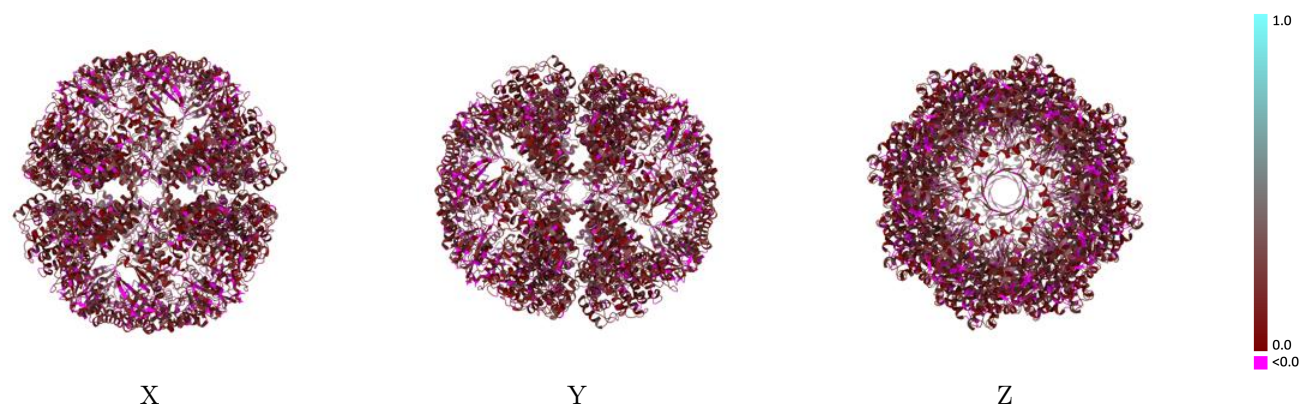
Y



Z

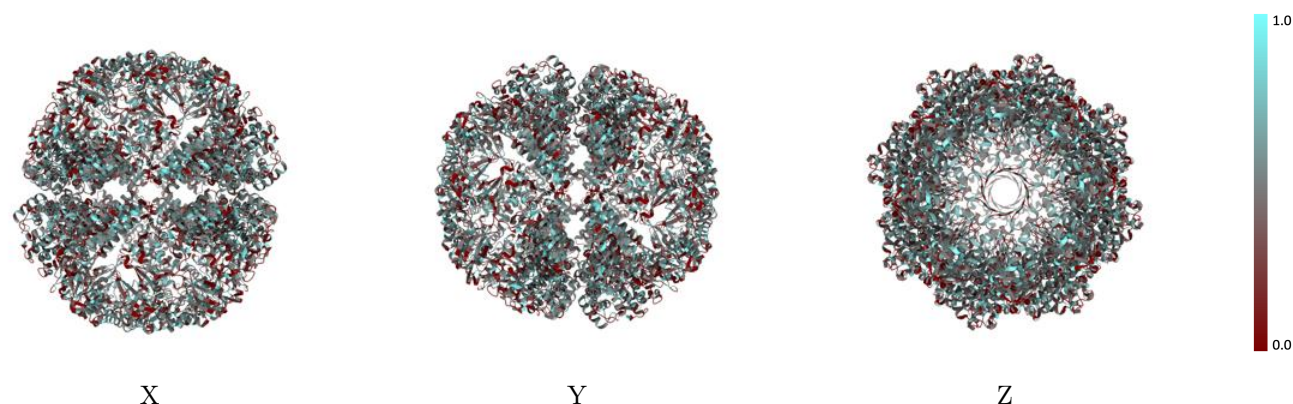
The images above show the 3D surface view of the map at the recommended contour level 0.28 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



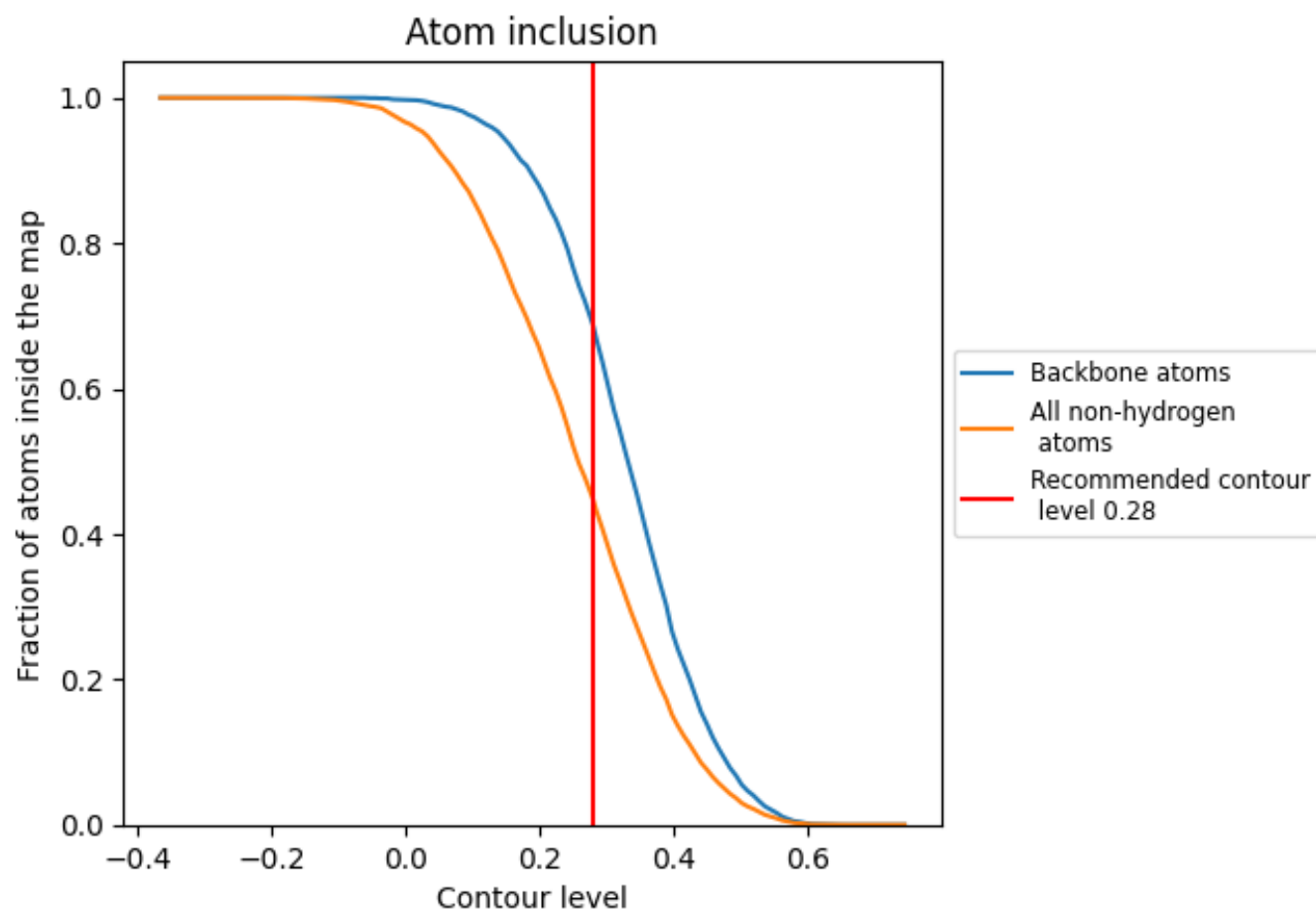
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.28).

9.4 Atom inclusion [i](#)



At the recommended contour level, 68% of all backbone atoms, 45% 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 (0.28) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.4457	<div></div> 0.1270
A	<div></div> 0.4465	<div></div> 0.1260
B	<div></div> 0.4457	<div></div> 0.1270
C	<div></div> 0.4462	<div></div> 0.1260
D	<div></div> 0.4459	<div></div> 0.1270
E	<div></div> 0.4462	<div></div> 0.1270
F	<div></div> 0.4457	<div></div> 0.1280
G	<div></div> 0.4462	<div></div> 0.1270
H	<div></div> 0.4457	<div></div> 0.1270
I	<div></div> 0.4451	<div></div> 0.1260
J	<div></div> 0.4454	<div></div> 0.1260
K	<div></div> 0.4457	<div></div> 0.1270
L	<div></div> 0.4454	<div></div> 0.1270
M	<div></div> 0.4457	<div></div> 0.1270
N	<div></div> 0.4457	<div></div> 0.1270
O	<div></div> 0.4457	<div></div> 0.1260
P	<div></div> 0.4451	<div></div> 0.1260

1.0

0.0

<0.0