



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 06:21 AM UTC

PDB ID : 1GW7 / pdb_00001gw7
EMDB ID : EMD-1011
Title : QUASI-ATOMIC RESOLUTION MODEL OF BACTERIOPHAGE PRD1 CAPSID, OBTAINED BY COMBINED CRYO-EM AND X-RAY CRYSTALLOGRAPHY.
Authors : San Martin, C.; Huiskonen, J.; Bamford, J.K.H.; Butcher, S.J.; Fuller, S.D.; Bamford, D.H.; Burnett, R.M.
Deposited on : 2002-03-08
Resolution : 13.50 Å(reported)
Based on initial model : 1GW7

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

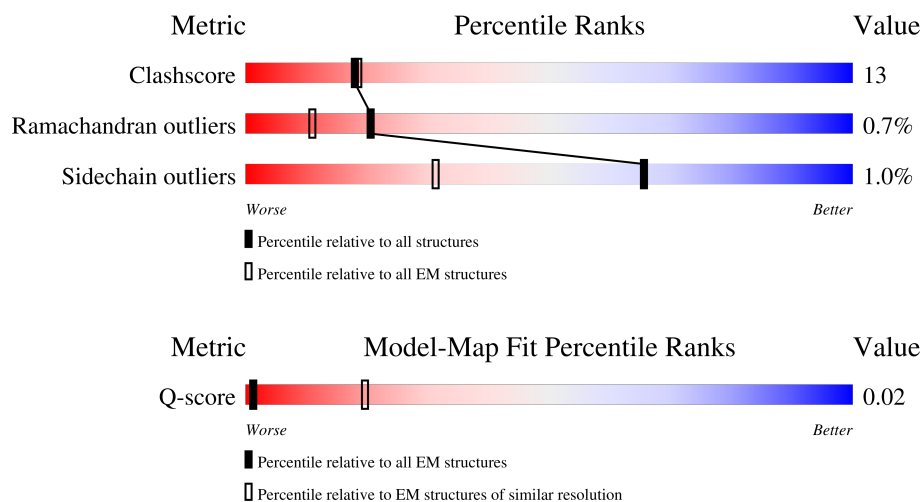
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 13.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	41 (15.50 - 16.50)

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	394	33% 76% 16% • 6%
1	B	394	34% 76% 16% • 5%
1	C	394	36% 78% 14% • 6%
1	D	394	37% 73% 19% • 6%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	E	394	41% 77% 15% 5%
1	F	394	39% 80% 12% 6%
1	G	394	42% 76% 16% 6%
1	H	394	37% 81% 12% 5%
1	I	394	37% 76% 16% 6%
1	J	394	35% 77% 16% 6%
1	K	394	34% 79% 13% 5%
1	L	394	28% 80% 13% 6%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 34170 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.

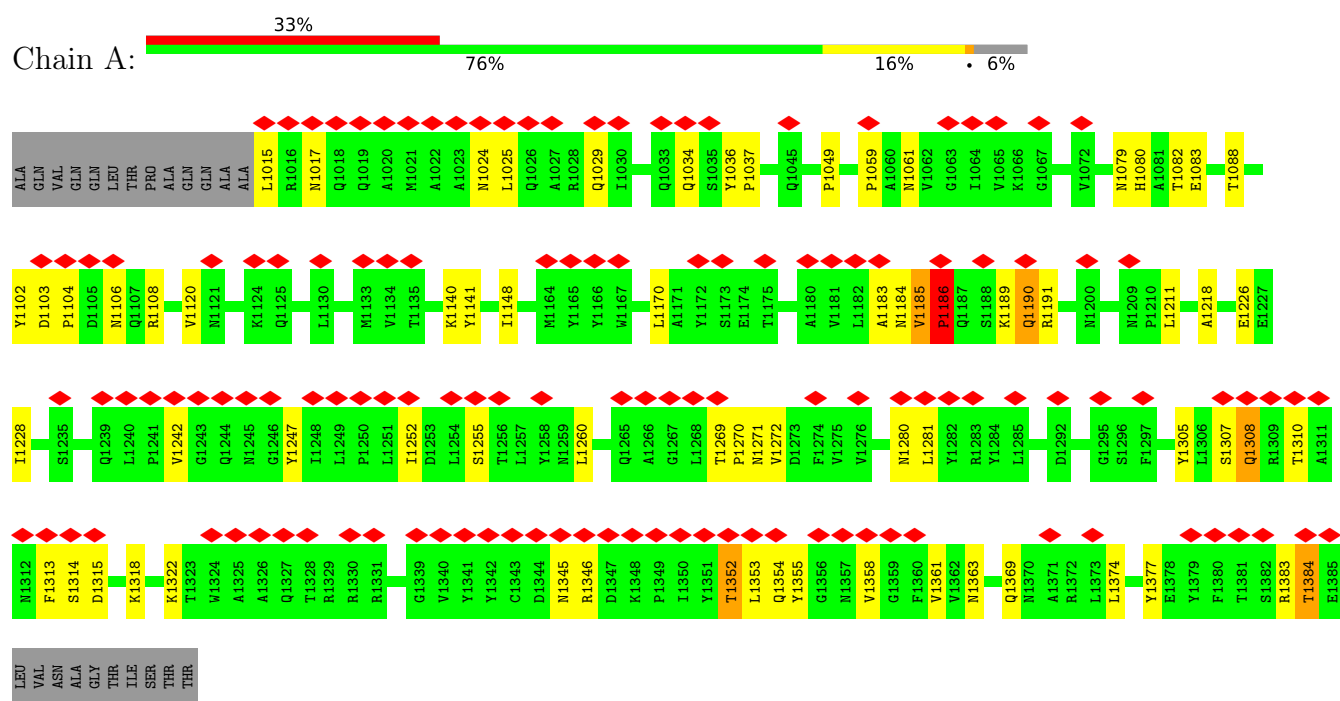
- Molecule 1 is a protein called MAJOR CAPSID PROTEIN.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	371	Total	C	N	O	S	0	1
			2837	1803	479	548	7		
1	B	373	Total	C	N	O	S	0	1
			2847	1804	485	551	7		
1	C	372	Total	C	N	O	S	0	1
			2861	1817	484	553	7		
1	D	371	Total	C	N	O	S	0	2
			2827	1795	478	547	7		
1	E	373	Total	C	N	O	S	0	1
			2847	1804	485	551	7		
1	F	372	Total	C	N	O	S	0	1
			2861	1817	484	553	7		
1	G	371	Total	C	N	O	S	0	1
			2837	1803	479	548	7		
1	H	373	Total	C	N	O	S	0	1
			2847	1804	485	551	7		
1	I	372	Total	C	N	O	S	0	1
			2861	1817	484	553	7		
1	J	371	Total	C	N	O	S	0	1
			2837	1803	479	548	7		
1	K	373	Total	C	N	O	S	0	1
			2847	1804	485	551	7		
1	L	372	Total	C	N	O	S	0	1
			2861	1817	484	553	7		

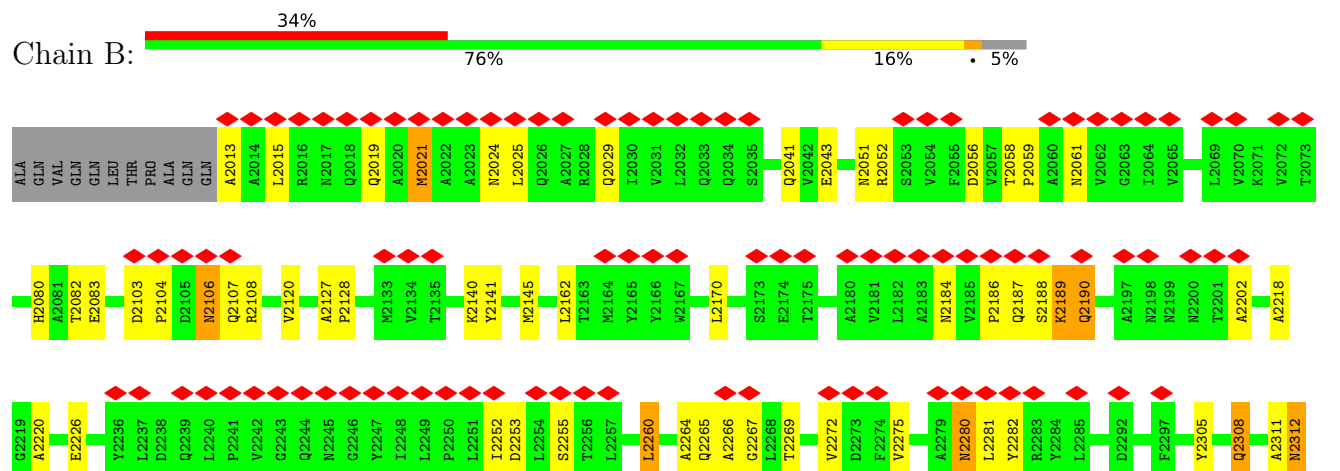
3 Residue-property plots

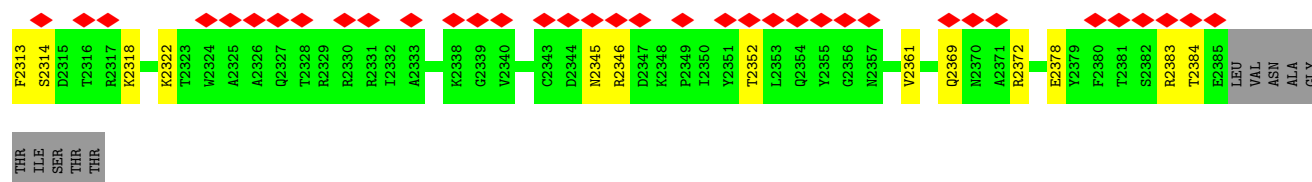
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: MAJOR CAPSID PROTEIN

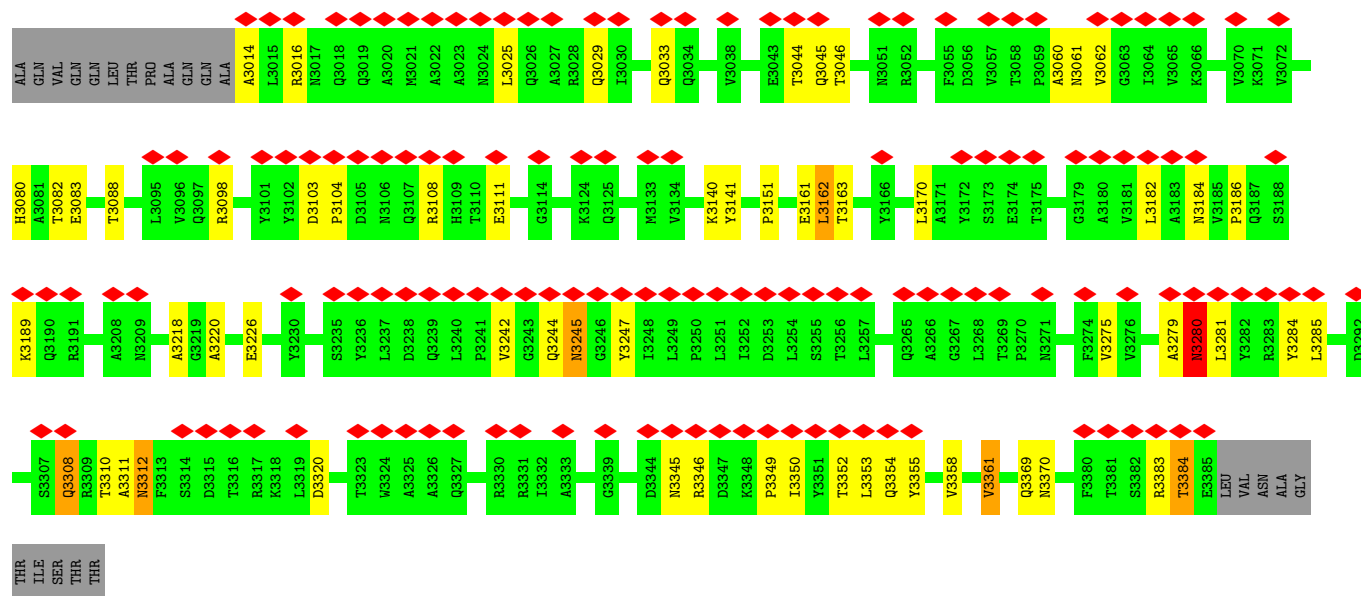
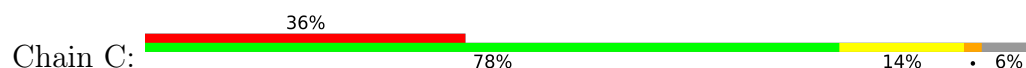


• Molecule 1: MAJOR CAPSID PROTEIN

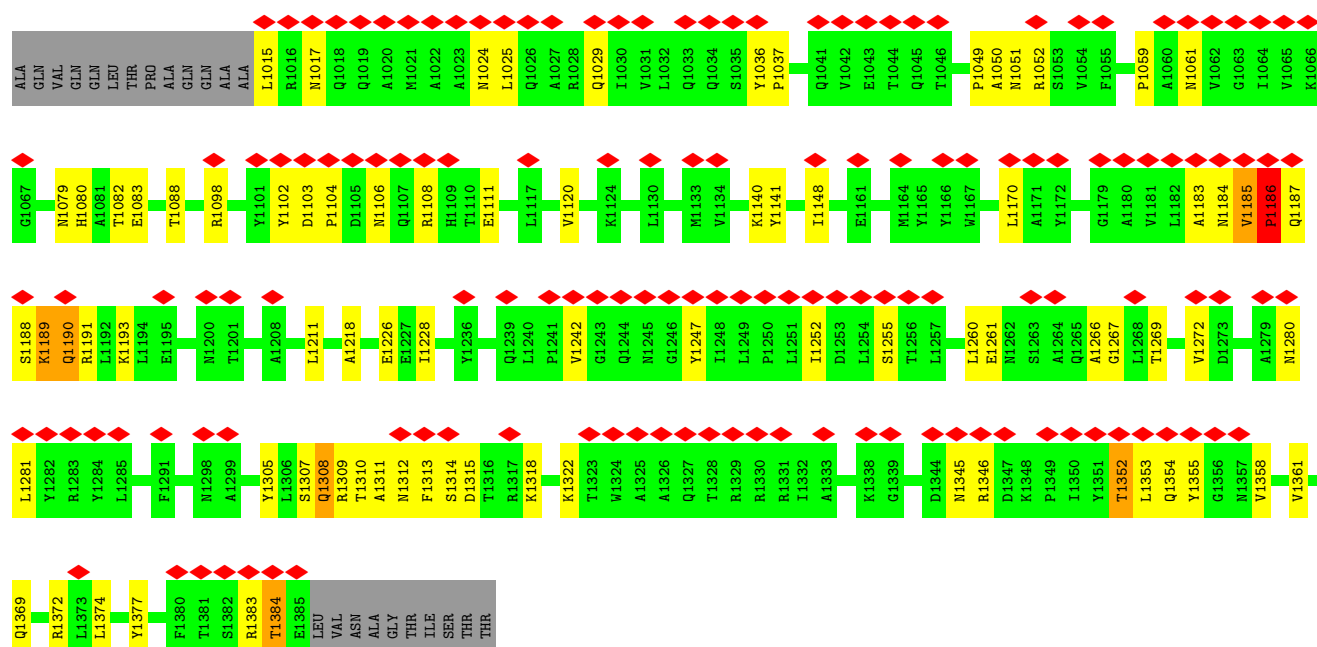




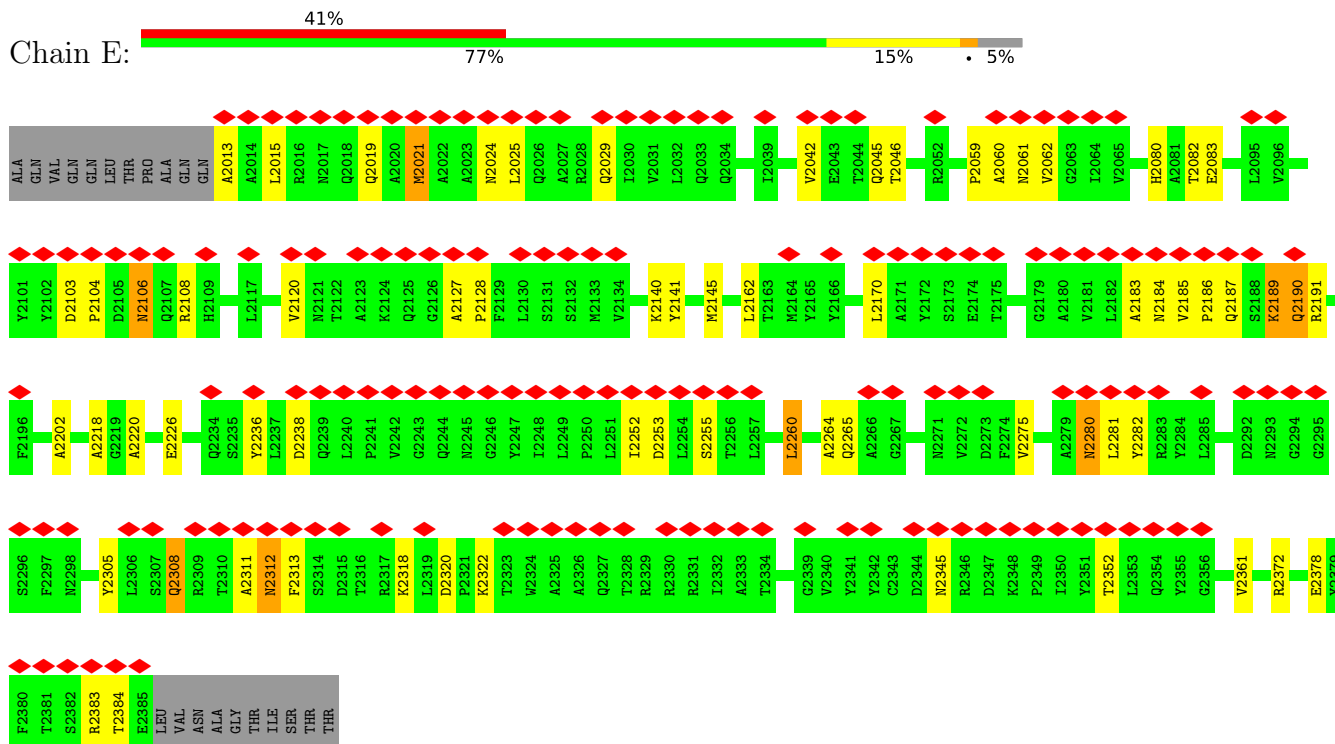
• Molecule 1: MAJOR CAPSID PROTEIN



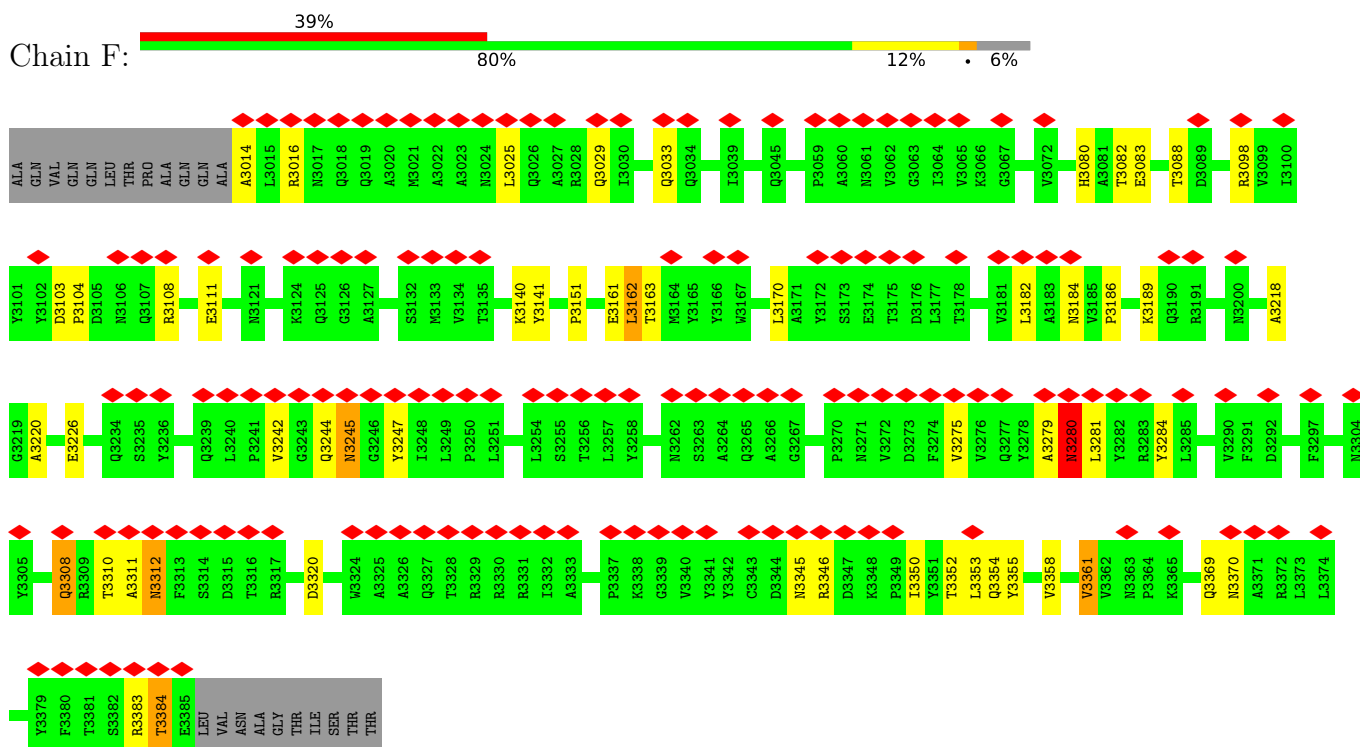
• Molecule 1: MAJOR CAPSID PROTEIN



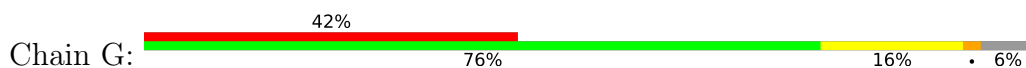
• Molecule 1: MAJOR CAPSID PROTEIN

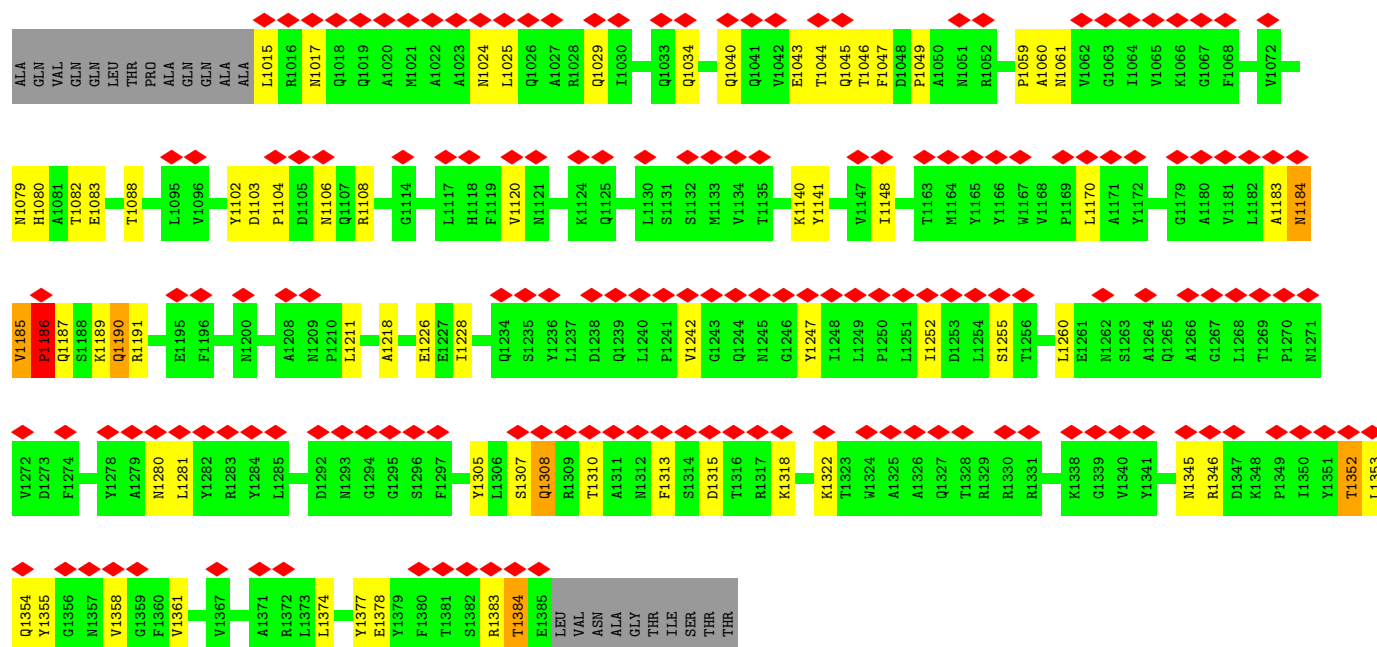


• Molecule 1: MAJOR CAPSID PROTEIN

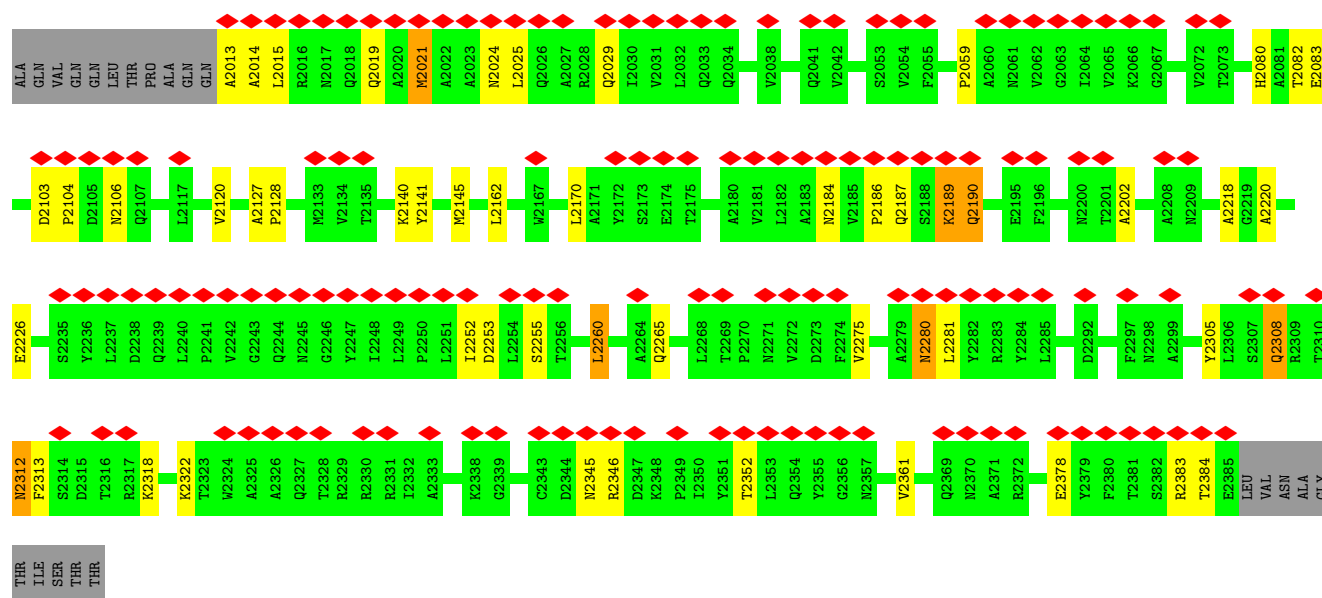
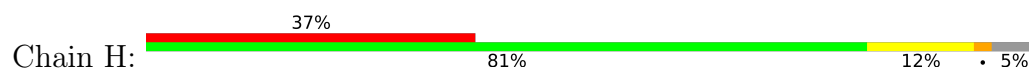


• Molecule 1: MAJOR CAPSID PROTEIN

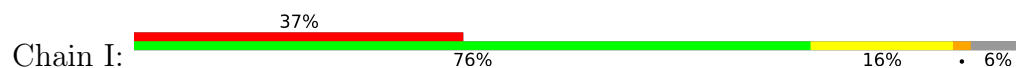


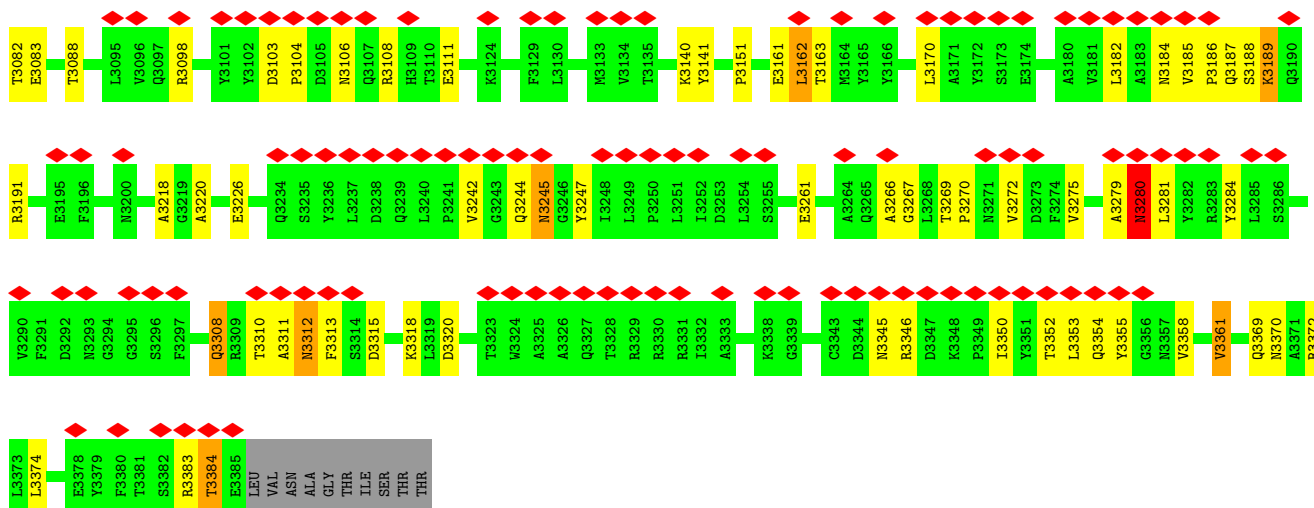


• Molecule 1: MAJOR CAPSID PROTEIN

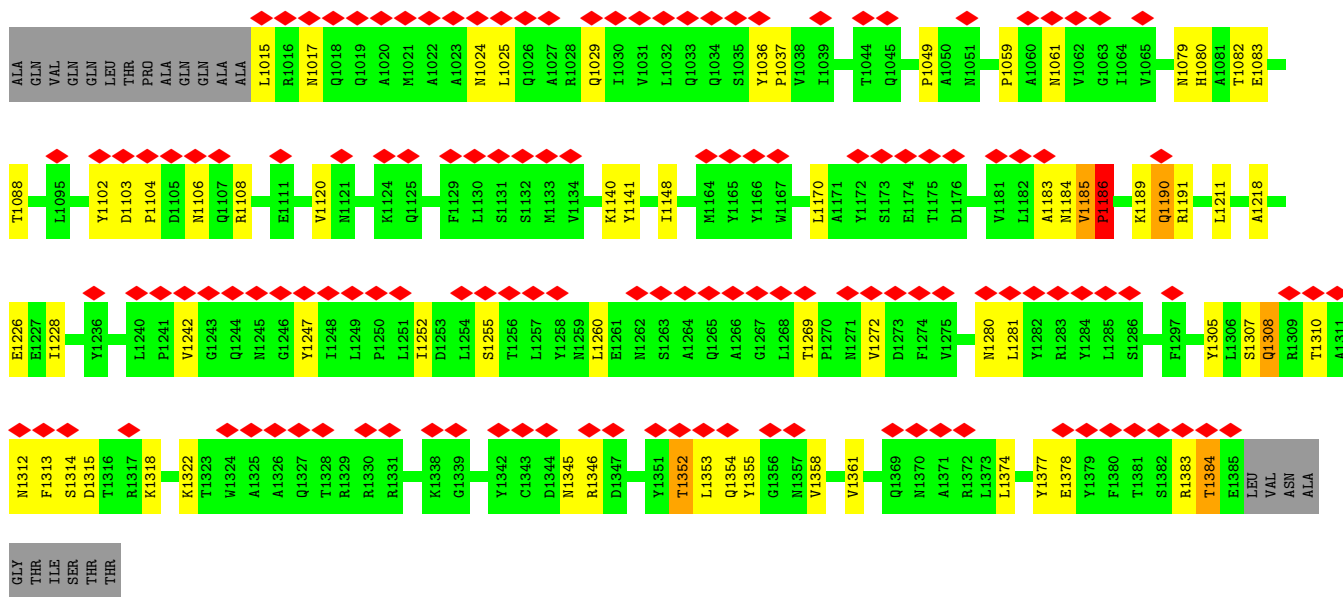
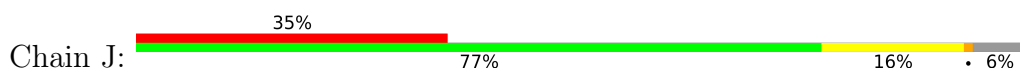


• Molecule 1: MAJOR CAPSID PROTEIN

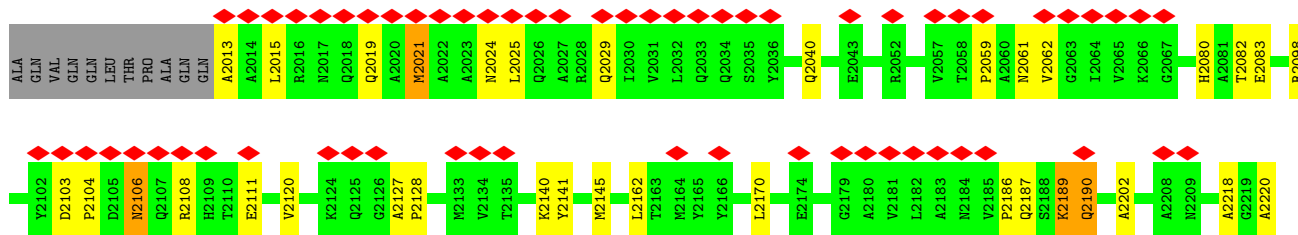
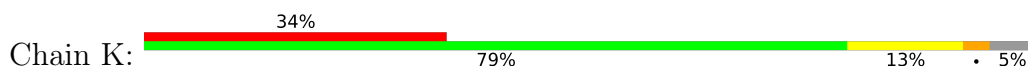


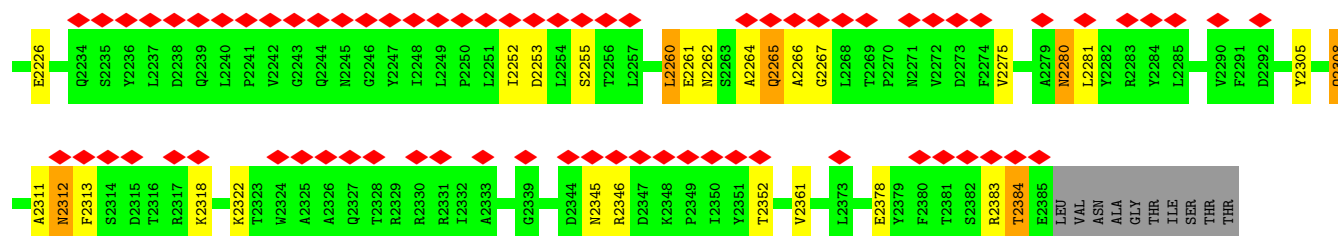


• Molecule 1: MAJOR CAPSID PROTEIN

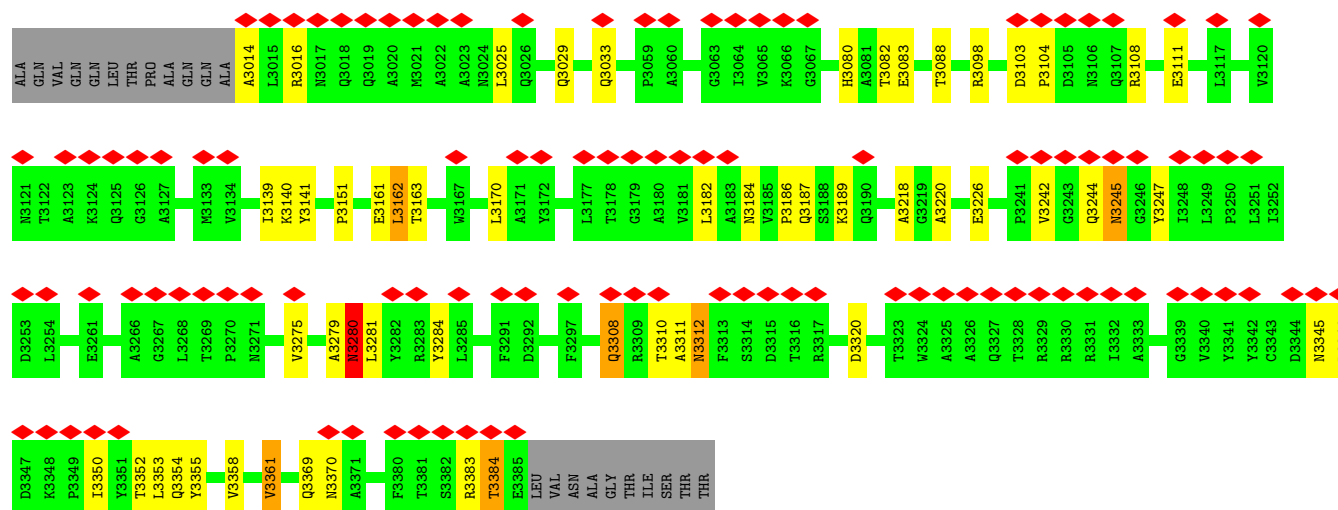
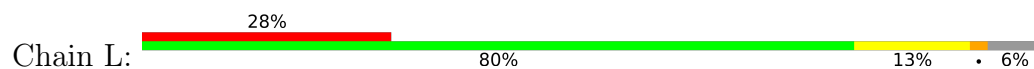


• Molecule 1: MAJOR CAPSID PROTEIN





• Molecule 1: MAJOR CAPSID PROTEIN



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	1468	Depositor
Resolution determination method	Not provided	
CTF correction method	PHASE RESTORATION BY CTF- MULTIPLICATION OF IMAGES. AMPLITUDE RESTORATION BY COMPARISON WITH QUASI- ATOMIC MODEL	Depositor
Microscope	FEI/PHILIPS CM200FEG	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	6	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	4100	Depositor
Magnification	45300	Depositor
Image detector	KODAK SO-163 FILM	Depositor
Maximum map value	27000.000	Depositor
Minimum map value	-16843.700	Depositor
Average map value	188.248	Depositor
Map value standard deviation	3602.720	Depositor
Recommended contour level	7770	Depositor
Map size ($\text{\AA}$)	880.64, 880.64, 880.64	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90, 90, 90	wwPDB
Pixel spacing ($\text{\AA}$)	3.44, 3.44, 3.44	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	0.48	1/2902 (0.0%)	1.01	14/3971 (0.4%)
1	B	0.51	1/2910 (0.0%)	0.99	12/3982 (0.3%)
1	C	0.49	1/2926 (0.0%)	1.00	13/4003 (0.3%)
1	D	0.48	1/2889 (0.0%)	1.00	13/3950 (0.3%)
1	E	0.51	1/2910 (0.0%)	0.99	12/3982 (0.3%)
1	F	0.49	1/2926 (0.0%)	1.00	13/4003 (0.3%)
1	G	0.49	1/2902 (0.0%)	1.00	14/3971 (0.4%)
1	H	0.52	1/2910 (0.0%)	0.99	12/3982 (0.3%)
1	I	0.49	1/2926 (0.0%)	1.00	13/4003 (0.3%)
1	J	0.48	1/2902 (0.0%)	1.01	13/3971 (0.3%)
1	K	0.52	1/2910 (0.0%)	0.99	12/3982 (0.3%)
1	L	0.49	1/2926 (0.0%)	1.00	14/4003 (0.3%)
All	All	0.50	12/34939 (0.0%)	1.00	155/47803 (0.3%)

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	2384	THR	C-N	-7.19	1.23	1.33
1	I	3384	THR	C-N	-7.17	1.23	1.33
1	D	1384	THR	C-N	-7.14	1.23	1.33
1	A	1384	THR	C-N	-7.12	1.23	1.33
1	H	2384	THR	C-N	-7.11	1.23	1.33

The worst 5 of 155 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	1170	LEU	N-CA-C	-10.90	97.52	112.12
1	D	1170	LEU	N-CA-C	-10.85	97.58	112.12
1	A	1170	LEU	N-CA-C	-10.81	97.64	112.12
1	G	1170	LEU	N-CA-C	-10.80	97.64	112.12
1	F	3170	LEU	N-CA-C	-10.54	97.99	112.12

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	2837	0	2709	79	0
1	B	2847	0	2723	241	0
1	C	2861	0	2751	76	0
1	D	2827	0	2695	237	0
1	E	2847	0	2724	143	0
1	F	2861	0	2752	35	0
1	G	2837	0	2710	127	0
1	H	2847	0	2729	38	0
1	I	2861	0	2740	185	0
1	J	2837	0	2711	69	0
1	K	2847	0	2729	92	0
1	L	2861	0	2752	35	0
All	All	34170	0	32725	892	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 13.

The worst 5 of 892 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:B:2264:ALA:CB	1:D:1272:VAL:CG1	1.74	1.61
1:B:2269:THR:CB	1:I:3372:ARG:HD3	1.28	1.60
1:B:2108:ARG:CD	1:D:1314:SER:HB2	1.36	1.54
1:B:2372:ARG:HD3	1:D:1269:THR:CG2	1.37	1.52
1:D:1189:LYS:CE	1:G:1060:ALA:HA	1.10	1.52

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	369/394 (94%)	349 (95%)	16 (4%)	4 (1%)	11	46
1	B	371/394 (94%)	349 (94%)	21 (6%)	1 (0%)	36	72
1	C	370/394 (94%)	348 (94%)	19 (5%)	3 (1%)	16	54
1	D	367/394 (93%)	347 (95%)	16 (4%)	4 (1%)	11	46
1	E	371/394 (94%)	349 (94%)	21 (6%)	1 (0%)	36	72
1	F	370/394 (94%)	348 (94%)	19 (5%)	3 (1%)	16	54
1	G	369/394 (94%)	349 (95%)	16 (4%)	4 (1%)	11	46
1	H	371/394 (94%)	349 (94%)	21 (6%)	1 (0%)	36	72
1	I	370/394 (94%)	348 (94%)	19 (5%)	3 (1%)	16	54
1	J	369/394 (94%)	349 (95%)	16 (4%)	4 (1%)	11	46
1	K	371/394 (94%)	349 (94%)	21 (6%)	1 (0%)	36	72
1	L	370/394 (94%)	348 (94%)	19 (5%)	3 (1%)	16	54
All	All	4438/4728 (94%)	4182 (94%)	224 (5%)	32 (1%)	20	56

5 of 32 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1185	VAL
1	C	3245	ASN
1	C	3280	ASN
1	C	3312	ASN
1	D	1185	VAL

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	292/325 (90%)	291 (100%)	1 (0%)	86	86
1	B	294/325 (90%)	288 (98%)	6 (2%)	48	66
1	C	298/325 (92%)	296 (99%)	2 (1%)	76	81
1	D	291/325 (90%)	290 (100%)	1 (0%)	86	86
1	E	294/325 (90%)	288 (98%)	6 (2%)	48	66
1	F	298/325 (92%)	296 (99%)	2 (1%)	76	81
1	G	292/325 (90%)	291 (100%)	1 (0%)	86	86
1	H	294/325 (90%)	288 (98%)	6 (2%)	48	66
1	I	298/325 (92%)	296 (99%)	2 (1%)	76	81
1	J	292/325 (90%)	291 (100%)	1 (0%)	86	86
1	K	294/325 (90%)	288 (98%)	6 (2%)	48	66
1	L	298/325 (92%)	296 (99%)	2 (1%)	76	81
All	All	3535/3900 (91%)	3499 (99%)	36 (1%)	65	78

5 of 36 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	K	2021	MET
1	L	3355	TYR
1	K	2106	ASN
1	K	2265	GLN
1	E	2162	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 128 such sidechains are listed below:

Mol	Chain	Res	Type
1	K	2265	GLN
1	K	2357	ASN
1	E	2107	GLN
1	E	2080	HIS
1	L	3079	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

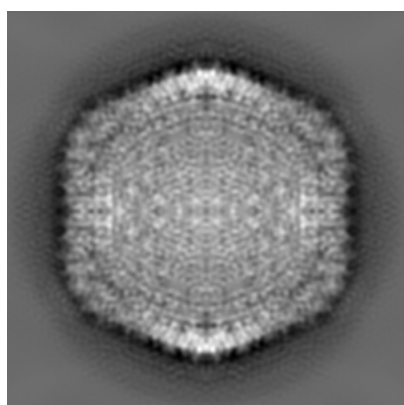
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-1011. These allow visual inspection of the internal detail of the map and identification of artifacts.

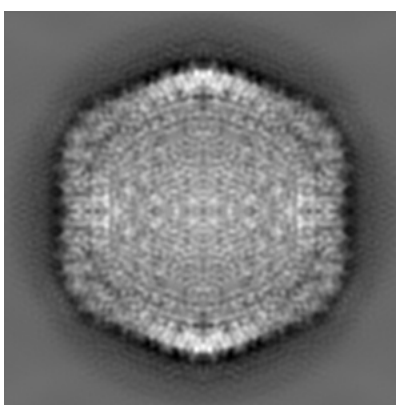
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

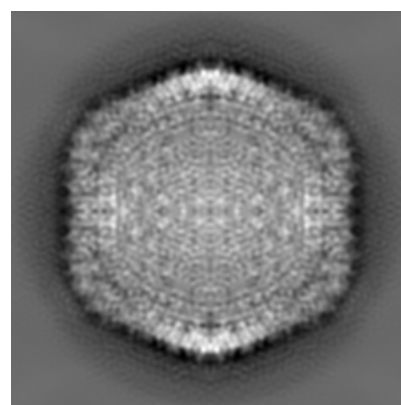
6.1.1 Primary map



X



Y

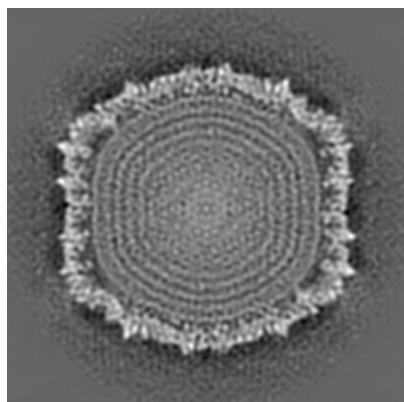


Z

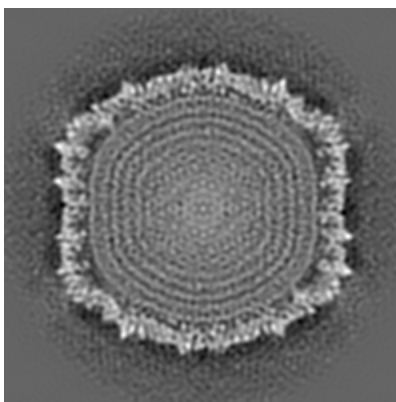
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

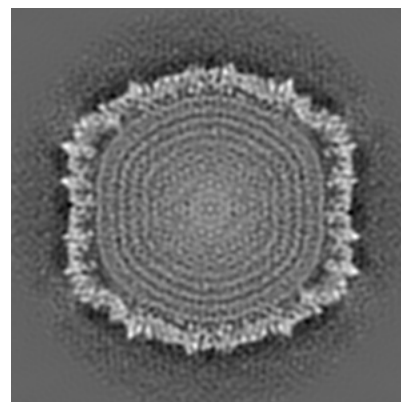
6.2.1 Primary map



X Index: 128



Y Index: 128

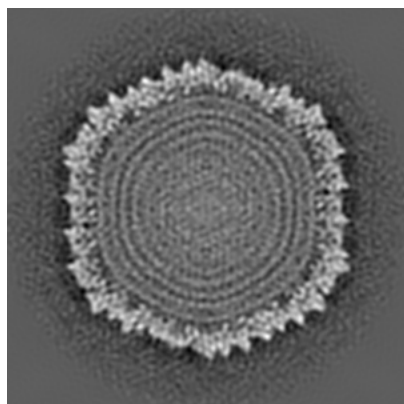


Z Index: 128

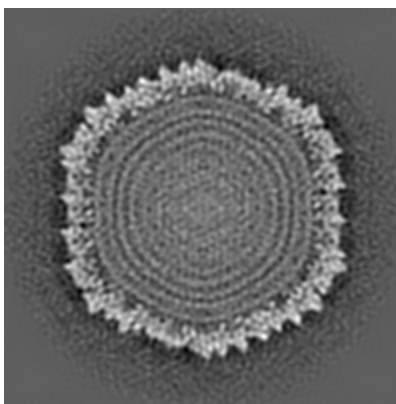
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

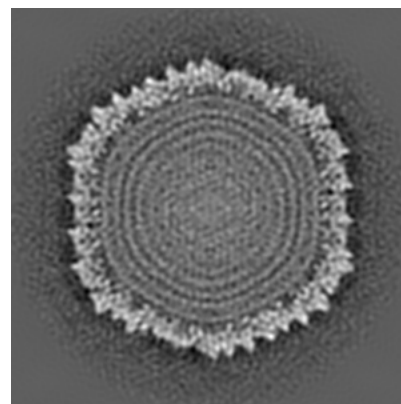
6.3.1 Primary map



X Index: 146



Y Index: 146

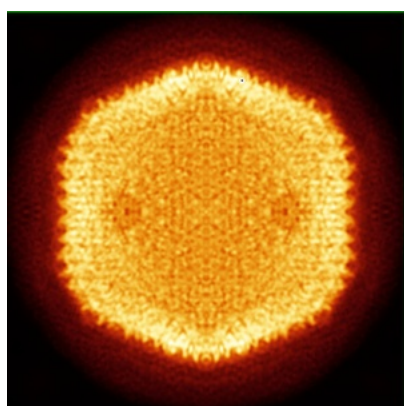


Z Index: 146

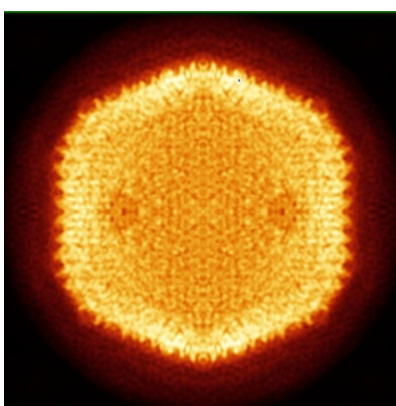
The images above show the largest variance slices of the map in three orthogonal directions.

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

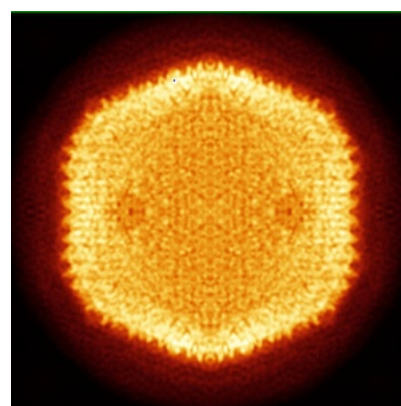
6.4.1 Primary map



X



Y

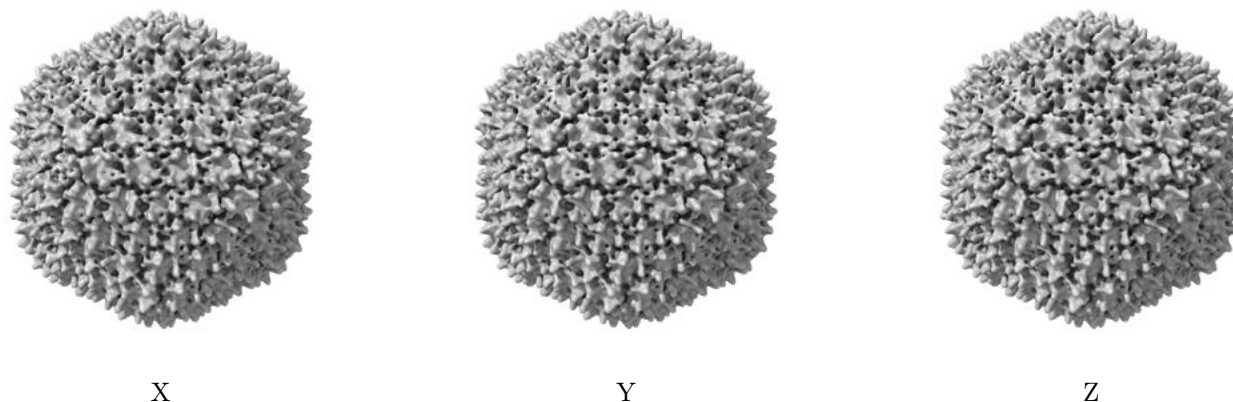


Z

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 7770.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

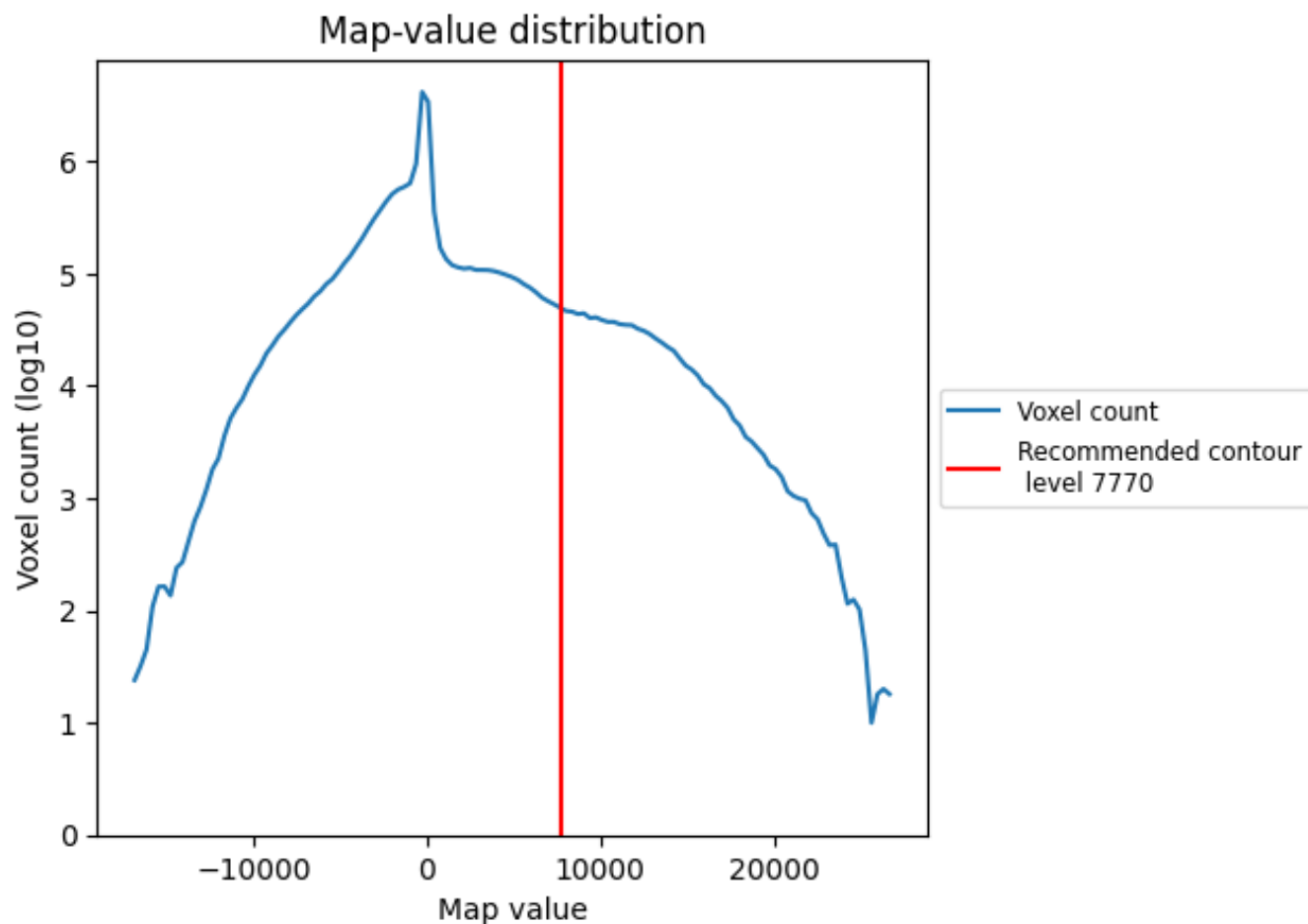
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

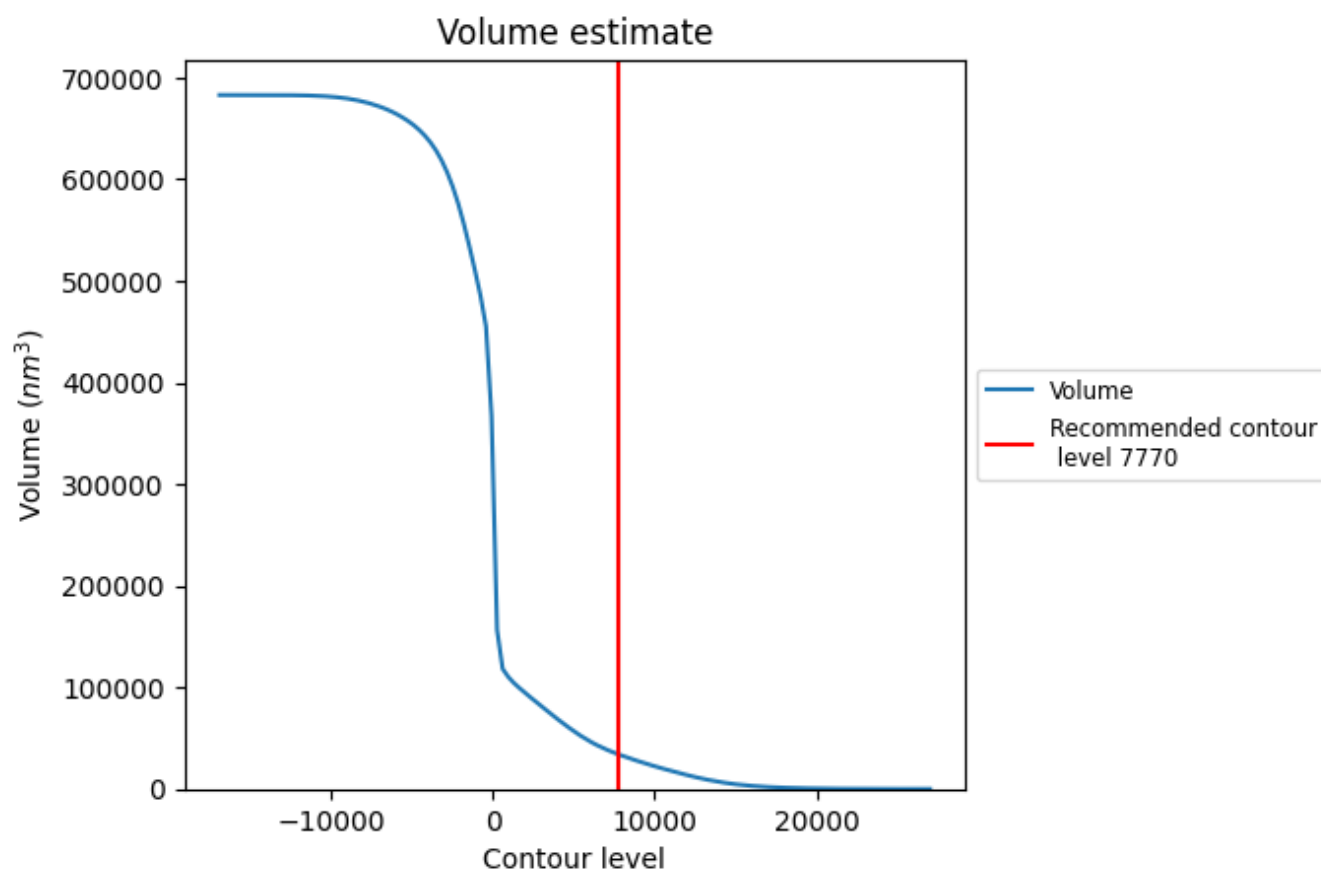
This section contains the results of statistical analysis of the map.

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.

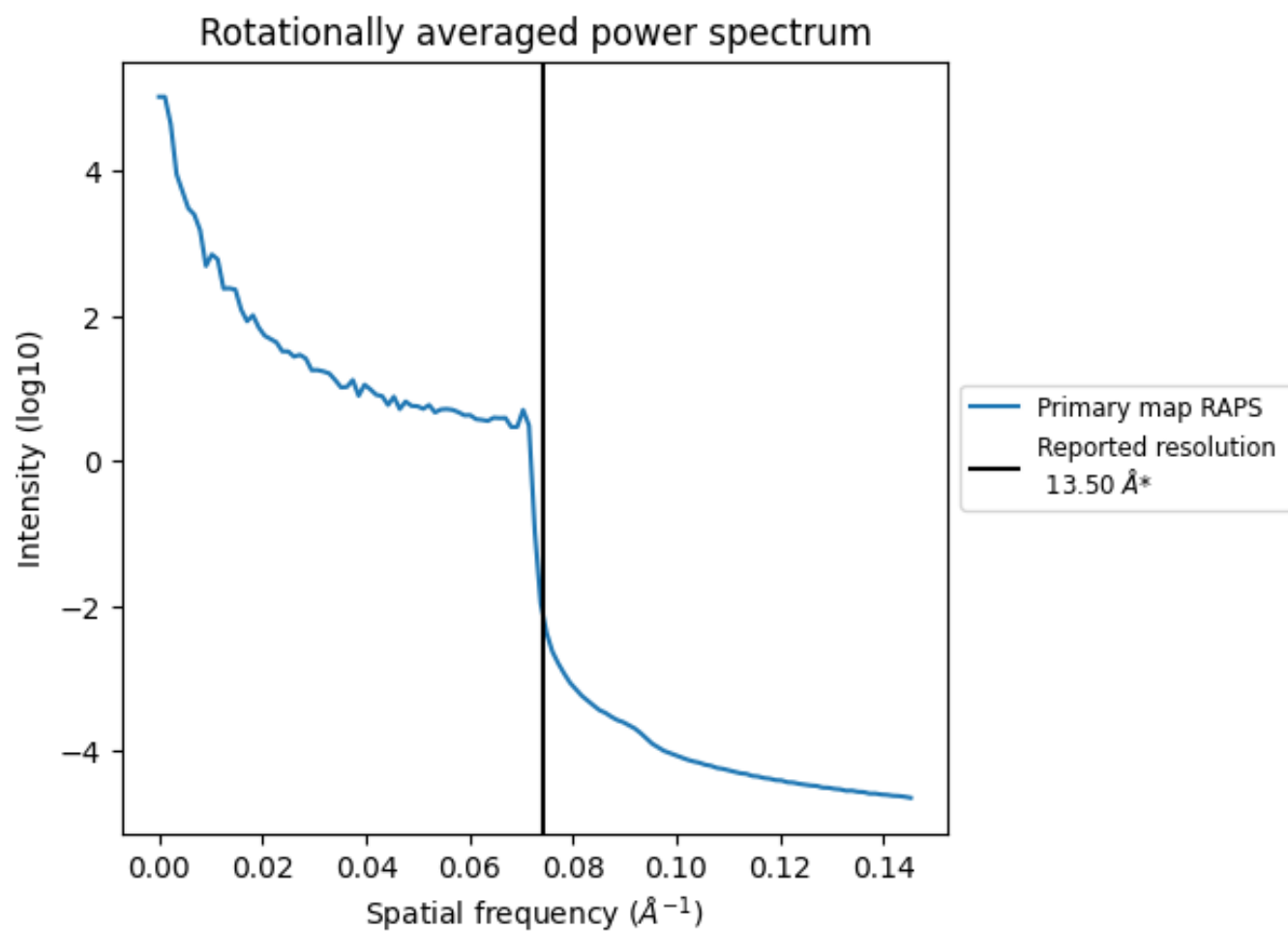
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 34042 nm^3 ; this corresponds to an approximate mass of 30751 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 0.074 Å⁻¹

8 Fourier-Shell correlation

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

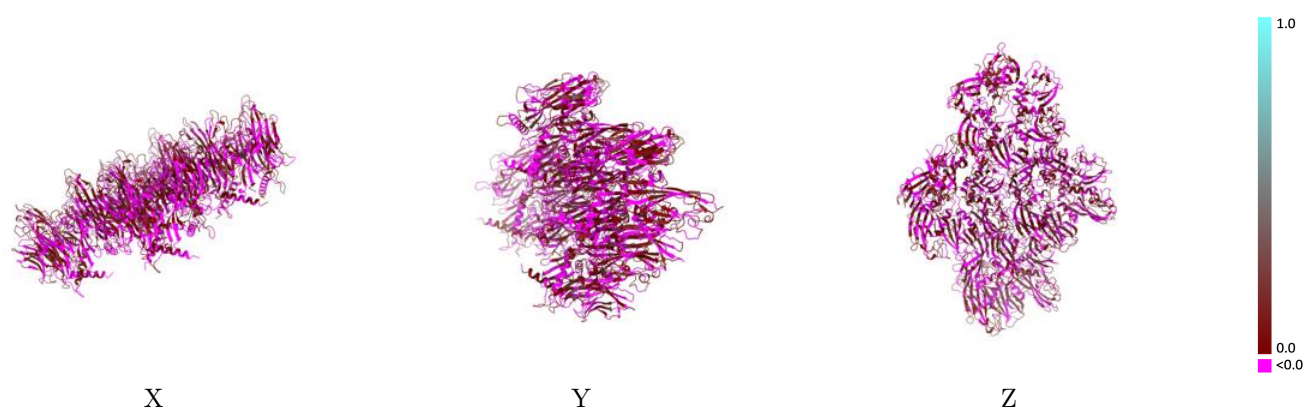
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-1011 and PDB model 1GW7. Per-residue inclusion information can be found in section [3](#) on page [5](#).

9.1 Map-model overlay [i](#)

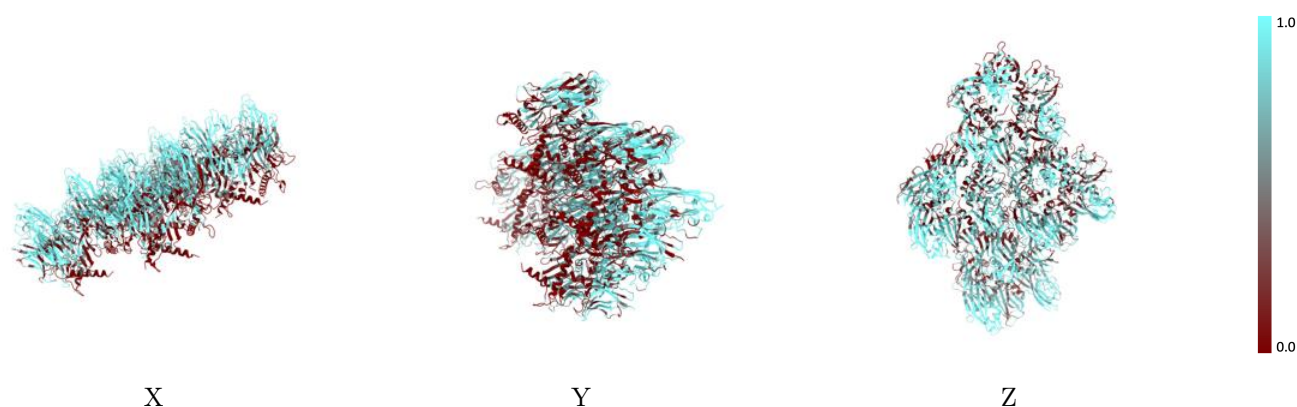
This section was not generated.

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



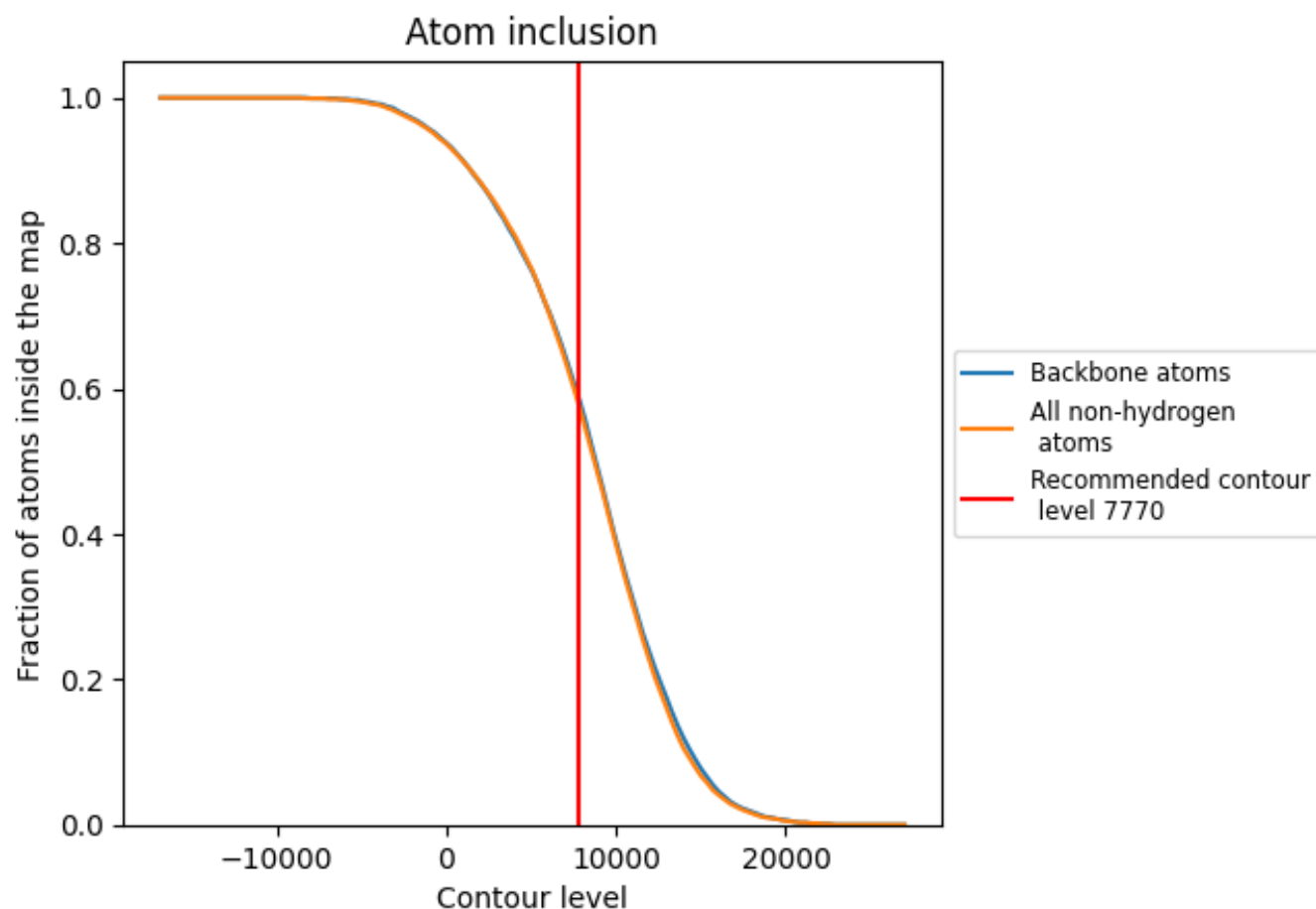
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (7770).

9.4 Atom inclusion [i](#)



At the recommended contour level, 59% of all backbone atoms, 58% 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 (7770) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.5810	0.0200
A	0.6070	0.0310
B	0.5930	0.0190
C	0.5800	0.0210
D	0.5730	0.0180
E	0.5410	0.0150
F	0.5560	0.0180
G	0.5250	0.0130
H	0.5640	0.0240
I	0.5710	0.0250
J	0.5900	0.0160
K	0.6070	0.0150
L	0.6620	0.0270

1.0

0.0

<0.0