Full wwPDB EM Validation Report

Dec 10, 2022 – 09:13 am GMT

PDB ID	:	5FJB
EMDB ID	:	EMD-3076
Title	:	Cyclophilin A Stabilize HIV-1 Capsid through a Novel Non- canonical Binding Site
Authors	:	Liu, C.; Perilla, J.R.; Ning, J.; Lu, M.; Hou, G.; Ramalhu, R.; Bedwell, G.J.; Ahn, J.; Shi, J.; Gronenborn, A.M.; Prevelige Jr, P.E.; Rousso, I.; Aiken, C.; Polenova, T.; Schulten, K.; Zhang, P.
Deposited on	:	2015-10-07
Resolution	:	9.00 Å(reported)
Based on initial model	:	3J4F

This is a Full wwPDB EM Validation 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 (i) 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 (i)) were used in the production of this report:

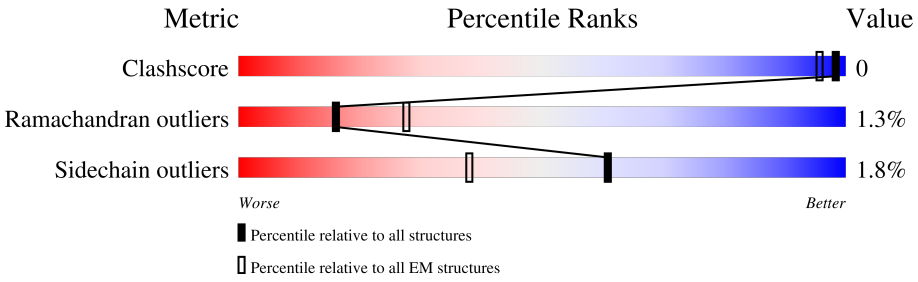
EMDB validation analysis	:	0.0.1.dev43
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 9.00 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
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 <=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	218	39% 79% 19% ·
1	B	218	37% 81% 18% ·
2	C	164	98% 84% 16% ·

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4664 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 GAG POLYPROTEIN.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	218	Total	C	N	O	S	0	0
			1703	1074	296	320	13		
1	B	218	Total	C	N	O	S	0	0
			1703	1074	296	320	13		

- Molecule 2 is a protein called PEPTIDYL-PROLYL CIS-TRANS ISOMERASE A.

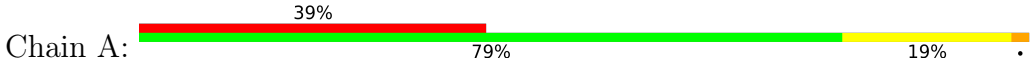
Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	164	Total	C	N	O	S	0	0
			1258	797	217	236	8		

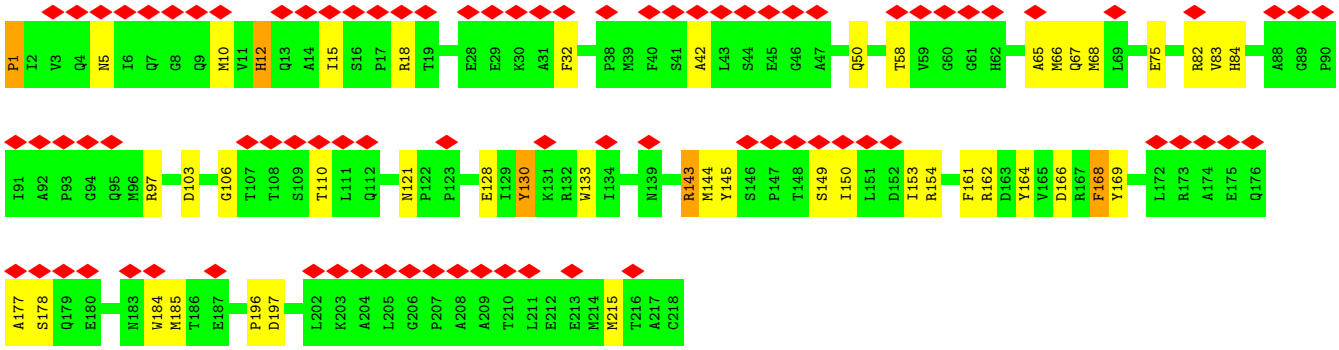
3 Residue-property plots (i)

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: GAG POLYPROTEIN

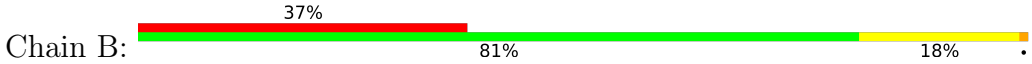
Chain A:

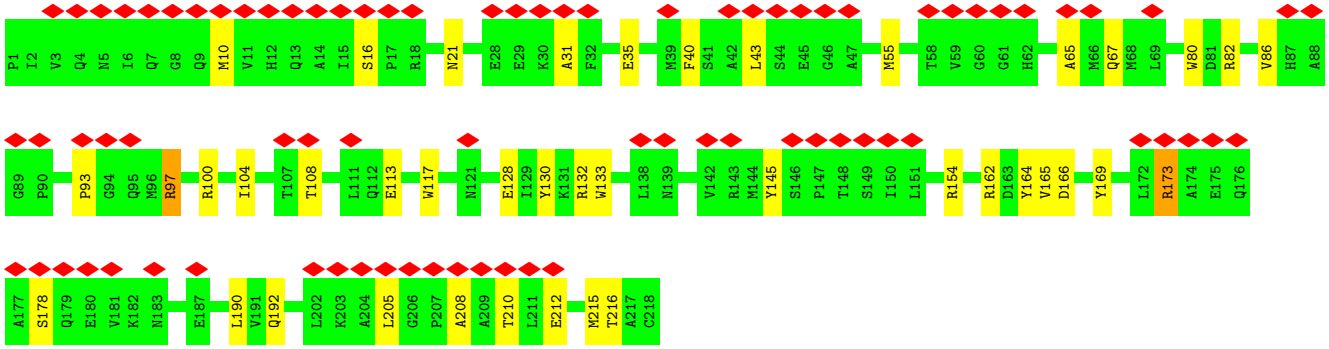




• Molecule 1: GAG POLYPROTEIN

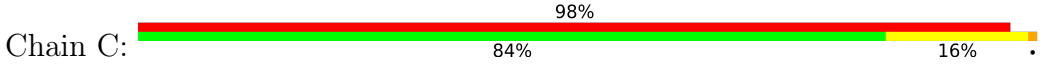
Chain B:

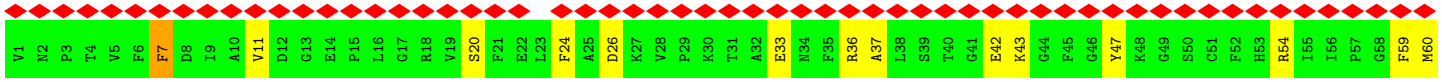


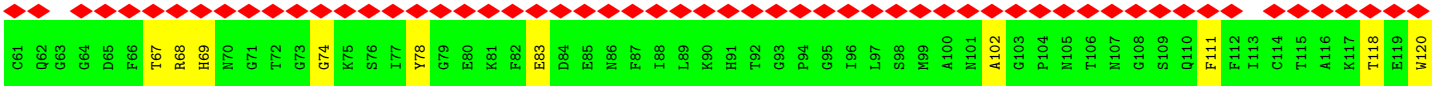


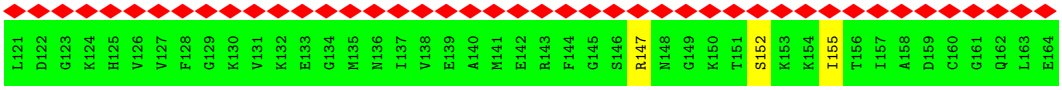
• Molecule 2: PEPTIDYL-PROLYL CIS-TRANS ISOMERASE A

Chain C:






C61 Q62 G63 G64 D65 F66 T67 R68 H69 N70 G71 T72 G73 G74 K75 S76 I77 Y78 G79 E80 K81 F82 E83 D84 E85 N86 F87 I88 L89 K90 H91 T92 G93 P94 G95 I96 L97 S98 M99 A100 N101 A102 G103 P104 N105 T106 N107 G108 S109 Q110 F111 F112 I113 C114 T115 A116 K117 T118 E119 W120


L121 D122 G123 K124 H125 V126 V127 F128 G129 K130 V131 K132 E133 G134 M135 N136 I137 V138 E139 A140 M141 E142 R143 F144 G145 S146 R147 N148 G149 K150 T151 S152 K153 K154 I155 T156 I157 A158 D159 C160 G161 Q162 L163 E164

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=Not provided°, rise=Not provided Å, axial sym=Not provided	Depositor
Number of segments used	Not provided	
Resolution determination method	Not provided	
CTF correction method	Not provided	
Microscope	FEI TECNAI F30	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	15	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	59000	Depositor
Image detector	KODAK SO-163 FILM	Depositor
Maximum map value	57.381	Depositor
Minimum map value	-44.915	Depositor
Average map value	0.244	Depositor
Map value standard deviation	11.385	Depositor
Recommended contour level	20.0	Depositor
Map size (Å)	462.15997, 462.15997, 275.59998	wwPDB
Map dimensions	436, 436, 260	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.06, 1.06, 1.06	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	1.66	10/1741 (0.6%)	1.99	32/2368 (1.4%)
1	B	1.59	9/1741 (0.5%)	1.95	35/2368 (1.5%)
2	C	1.67	8/1286 (0.6%)	1.86	22/1723 (1.3%)
All	All	1.64	27/4768 (0.6%)	1.94	89/6459 (1.4%)

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	7
1	B	0	2
2	C	0	2
All	All	0	11

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	106	GLY	CA-C	-7.34	1.40	1.51
1	B	16	SER	CA-CB	6.92	1.63	1.52
2	C	83	GLU	CD-OE1	6.88	1.33	1.25
1	B	65	ALA	N-CA	-6.44	1.33	1.46
1	A	145	TYR	CE1-CZ	6.38	1.46	1.38
1	B	205	LEU	C-N	6.19	1.44	1.33
1	B	31	ALA	N-CA	6.05	1.58	1.46
1	A	50	GLN	CA-CB	6.04	1.67	1.53
2	C	36	ARG	CD-NE	5.60	1.55	1.46
2	C	47	TYR	CG-CD2	5.50	1.46	1.39
1	B	133	TRP	CA-CB	5.43	1.66	1.53
2	C	43	LYS	CA-CB	5.40	1.65	1.53
1	A	133	TRP	CE3-CZ3	5.38	1.47	1.38
2	C	147	ARG	CZ-NH1	-5.37	1.26	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	74	GLY	N-CA	5.37	1.54	1.46
1	B	113	GLU	CB-CG	5.31	1.62	1.52
1	A	12	HIS	CG-CD2	-5.28	1.26	1.35
1	A	178	SER	CA-CB	5.26	1.60	1.52
2	C	47	TYR	CZ-OH	5.19	1.46	1.37
1	A	133	TRP	NE1-CE2	-5.18	1.30	1.37
1	B	128	GLU	CD-OE1	5.18	1.31	1.25
2	C	20	SER	CA-CB	5.17	1.60	1.52
1	A	168	PHE	CG-CD1	5.17	1.46	1.38
1	B	178	SER	C-N	5.14	1.45	1.34
1	A	75	GLU	CB-CG	5.11	1.61	1.52
1	A	168	PHE	CE1-CZ	5.08	1.47	1.37
1	B	190	LEU	CA-CB	5.08	1.65	1.53

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	154	ARG	NE-CZ-NH1	22.44	131.52	120.30
1	B	97	ARG	NE-CZ-NH2	-20.21	110.19	120.30
1	A	82	ARG	NE-CZ-NH1	18.10	129.35	120.30
1	B	132	ARG	NE-CZ-NH1	15.42	128.01	120.30
1	B	97	ARG	NE-CZ-NH1	15.29	127.95	120.30
2	C	36	ARG	NE-CZ-NH1	14.37	127.49	120.30
1	A	82	ARG	NE-CZ-NH2	-11.06	114.77	120.30
1	B	132	ARG	NE-CZ-NH2	-10.50	115.05	120.30
1	A	154	ARG	NE-CZ-NH2	-10.29	115.15	120.30
1	A	162	ARG	NE-CZ-NH1	10.07	125.34	120.30
1	B	145	TYR	CB-CG-CD2	-10.00	115.00	121.00
1	B	100	ARG	NE-CZ-NH1	9.77	125.18	120.30
1	B	173	ARG	NE-CZ-NH2	-9.66	115.47	120.30
1	B	162	ARG	NE-CZ-NH1	9.35	124.97	120.30
1	B	169	TYR	CD1-CE1-CZ	-9.34	111.40	119.80
1	A	143	ARG	NE-CZ-NH2	-9.01	115.79	120.30
2	C	36	ARG	NH1-CZ-NH2	-8.93	109.58	119.40
1	A	168	PHE	CB-CG-CD1	8.70	126.89	120.80
1	A	185	MET	CG-SD-CE	-8.68	86.31	100.20
1	A	110	THR	CA-CB-CG2	-8.17	100.96	112.40
2	C	68	ARG	NE-CZ-NH1	8.15	124.37	120.30
1	A	197	ASP	CB-CG-OD1	7.96	125.46	118.30
1	A	130	TYR	CB-CG-CD2	-7.61	116.43	121.00
2	C	152	SER	CB-CA-C	-7.52	95.81	110.10
1	B	164	TYR	CB-CG-CD2	-7.40	116.56	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	164	TYR	CB-CG-CD2	7.37	125.42	121.00
1	A	168	PHE	CB-CG-CD2	-7.30	115.69	120.80
1	A	66	MET	CG-SD-CE	-7.29	88.53	100.20
2	C	59	PHE	CB-CG-CD1	-7.29	115.70	120.80
1	B	169	TYR	CZ-CE2-CD2	-7.19	113.33	119.80
1	B	166	ASP	CB-CG-OD1	7.04	124.64	118.30
2	C	37	ALA	O-C-N	-6.99	111.52	122.70
1	A	97	ARG	NE-CZ-NH1	6.99	123.79	120.30
2	C	120	TRP	CB-CG-CD2	6.99	135.68	126.60
1	B	215	MET	CG-SD-CE	-6.86	89.22	100.20
1	A	144	MET	CG-SD-CE	-6.85	89.24	100.20
2	C	102	ALA	N-CA-CB	-6.80	100.59	110.10
2	C	78	TYR	CB-CG-CD1	-6.78	116.94	121.00
1	A	83	VAL	CG1-CB-CG2	-6.74	100.12	110.90
1	A	215	MET	CA-CB-CG	6.64	124.58	113.30
2	C	67	THR	N-CA-CB	6.42	122.49	110.30
1	B	145	TYR	CB-CG-CD1	6.39	124.84	121.00
2	C	42	GLU	OE1-CD-OE2	-6.28	115.76	123.30
1	B	166	ASP	CB-CG-OD2	-6.22	112.70	118.30
1	A	166	ASP	CB-CG-OD1	6.22	123.90	118.30
1	B	55	MET	CG-SD-CE	-6.12	90.41	100.20
1	B	10	MET	N-CA-C	6.08	127.40	111.00
2	C	59	PHE	CG-CD1-CE1	-5.97	114.23	120.80
2	C	33	GLU	OE1-CD-OE2	-5.86	116.27	123.30
2	C	24	PHE	CB-CG-CD1	5.79	124.85	120.80
1	A	128	GLU	OE1-CD-OE2	-5.78	116.37	123.30
1	B	169	TYR	CG-CD2-CE2	5.77	125.92	121.30
1	A	161	PHE	CB-CG-CD1	5.74	124.81	120.80
1	B	212	GLU	OE1-CD-OE2	-5.67	116.50	123.30
1	B	21	ASN	O-C-N	-5.65	113.66	122.70
1	B	108	THR	CA-CB-CG2	-5.62	104.54	112.40
1	A	68	MET	CG-SD-CE	-5.60	91.25	100.20
1	A	154	ARG	NH1-CZ-NH2	-5.52	113.33	119.40
1	A	103	ASP	CB-CG-OD2	5.47	123.22	118.30
2	C	118	THR	OG1-CB-CG2	-5.47	97.42	110.00
2	C	7	PHE	CB-CA-C	-5.46	99.49	110.40
1	A	84	HIS	CA-CB-CG	-5.43	104.37	113.60
1	B	169	TYR	CB-CG-CD1	5.40	124.24	121.00
1	B	104	ILE	O-C-N	-5.38	114.09	122.70
1	B	82	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	B	40	PHE	CB-CG-CD1	-5.33	117.07	120.80
1	B	93	PRO	N-CA-C	5.32	125.92	112.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	152	SER	N-CA-CB	5.30	118.46	110.50
2	C	26	ASP	CB-CG-OD1	5.30	123.07	118.30
1	A	130	TYR	CG-CD1-CE1	-5.29	117.06	121.30
1	B	208	ALA	CB-CA-C	5.28	118.03	110.10
1	B	67	GLN	CG-CD-OE1	5.28	132.16	121.60
1	A	42	ALA	N-CA-CB	-5.26	102.73	110.10
2	C	54	ARG	CD-NE-CZ	5.26	130.96	123.60
2	C	7	PHE	N-CA-CB	5.25	120.05	110.60
1	A	15	ILE	O-C-N	5.25	131.09	122.70
1	A	65	ALA	O-C-N	-5.21	114.36	122.70
1	B	130	TYR	CG-CD2-CE2	-5.20	117.14	121.30
1	A	133	TRP	NE1-CE2-CD2	5.19	112.49	107.30
1	B	210	THR	N-CA-CB	5.17	120.13	110.30
1	B	117	TRP	NE1-CE2-CZ2	5.17	136.09	130.40
1	B	205	LEU	CB-CG-CD2	5.17	119.79	111.00
2	C	147	ARG	NE-CZ-NH1	-5.16	117.72	120.30
1	B	43	LEU	CB-CG-CD2	5.15	119.76	111.00
1	A	184	TRP	CB-CG-CD1	-5.04	120.45	127.00
1	B	165	VAL	CA-CB-CG2	-5.03	103.36	110.90
1	A	103	ASP	OD1-CG-OD2	-5.01	113.78	123.30
1	B	80	TRP	CB-CG-CD2	-5.00	120.09	126.60
2	C	78	TYR	CB-CG-CD2	5.00	124.00	121.00

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	PRO	Peptide
1	A	121	ASN	Peptide
1	A	130	TYR	Sidechain
1	A	168	PHE	Sidechain
1	A	169	TYR	Sidechain
1	A	18	ARG	Sidechain
1	A	32	PHE	Sidechain
1	B	173	ARG	Sidechain
1	B	97	ARG	Sidechain
2	C	111	PHE	Sidechain
2	C	7	PHE	Sidechain

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	1703	0	1704	1	0
1	B	1703	0	1704	0	0
2	C	1258	0	1228	1	0
All	All	4664	0	4636	2	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 0.

All (2) 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:A:150:ILE:HD12	1:A:153:ILE:HG13	1.82	0.61
2:C:11:VAL:HG22	2:C:155:ILE:HD12	2.03	0.41

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	216/218 (99%)	199 (92%)	12 (6%)	5 (2%)	6 34
1	B	216/218 (99%)	205 (95%)	9 (4%)	2 (1%)	17 57
2	C	162/164 (99%)	149 (92%)	12 (7%)	1 (1%)	25 66
All	All	594/600 (99%)	553 (93%)	33 (6%)	8 (1%)	16 48

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	MET
1	A	149	SER
1	A	177	ALA
2	C	69	HIS
1	A	5	ASN
1	B	35	GLU
1	B	86	VAL
1	A	196	PRO

5.3.2 Protein sidechains (i)

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	186/186 (100%)	181 (97%)	5 (3%)	44	65
1	B	186/186 (100%)	183 (98%)	3 (2%)	62	79
2	C	132/132 (100%)	131 (99%)	1 (1%)	81	89
All	All	504/504 (100%)	495 (98%)	9 (2%)	61	77

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	PRO
1	A	12	HIS
1	A	58	THR
1	A	67	GLN
1	A	143	ARG
1	B	154	ARG
1	B	192	GLN
1	B	216	THR
2	C	60	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	87	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

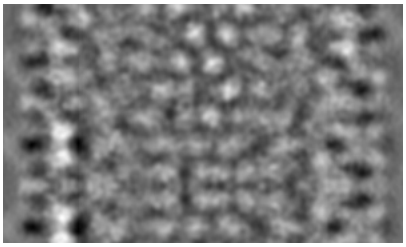
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-3076. These allow visual inspection of the internal detail of the map and identification of artifacts.

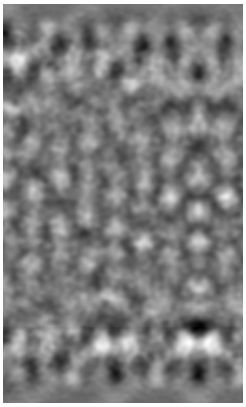
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

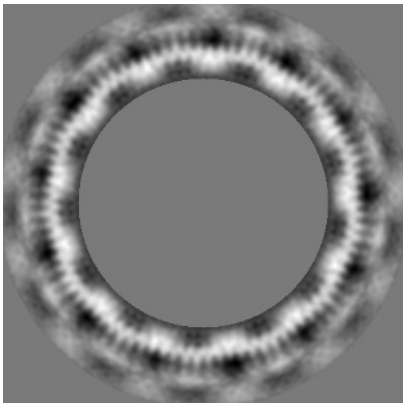
6.1.1 Primary map



X



Y



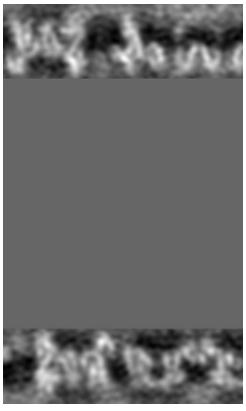
Z

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

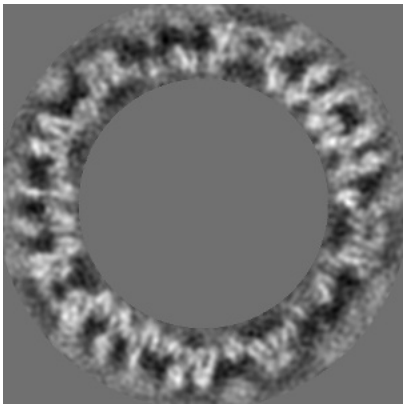
6.2 Central slices

6.2.1 Primary map

X Index: 218



Y Index: 218

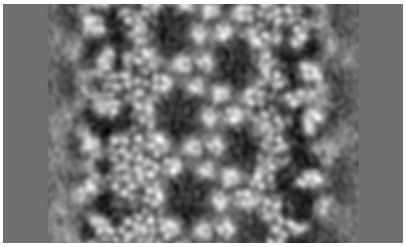


Z Index: 130

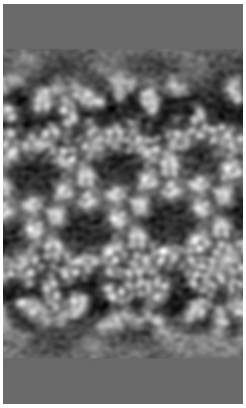
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

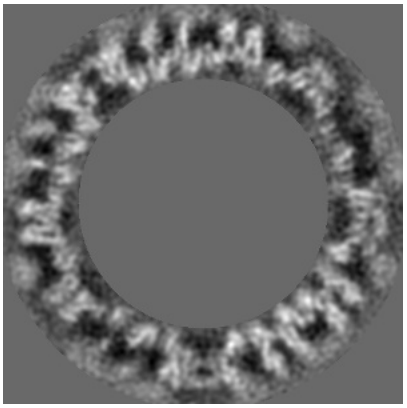
6.3.1 Primary map



X Index: 366



Y Index: 367

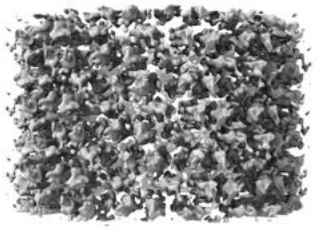


Z Index: 56

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

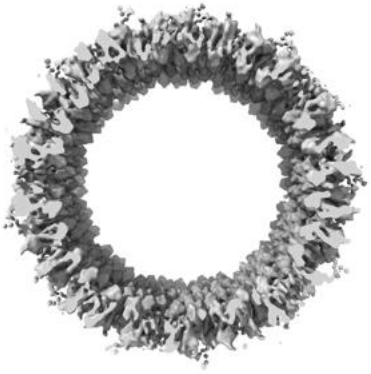
6.4 Orthogonal surface views (i)

6.4.1 Primary map



X

Y



Z

The images above show the 3D surface view of the map at the recommended contour level 20.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

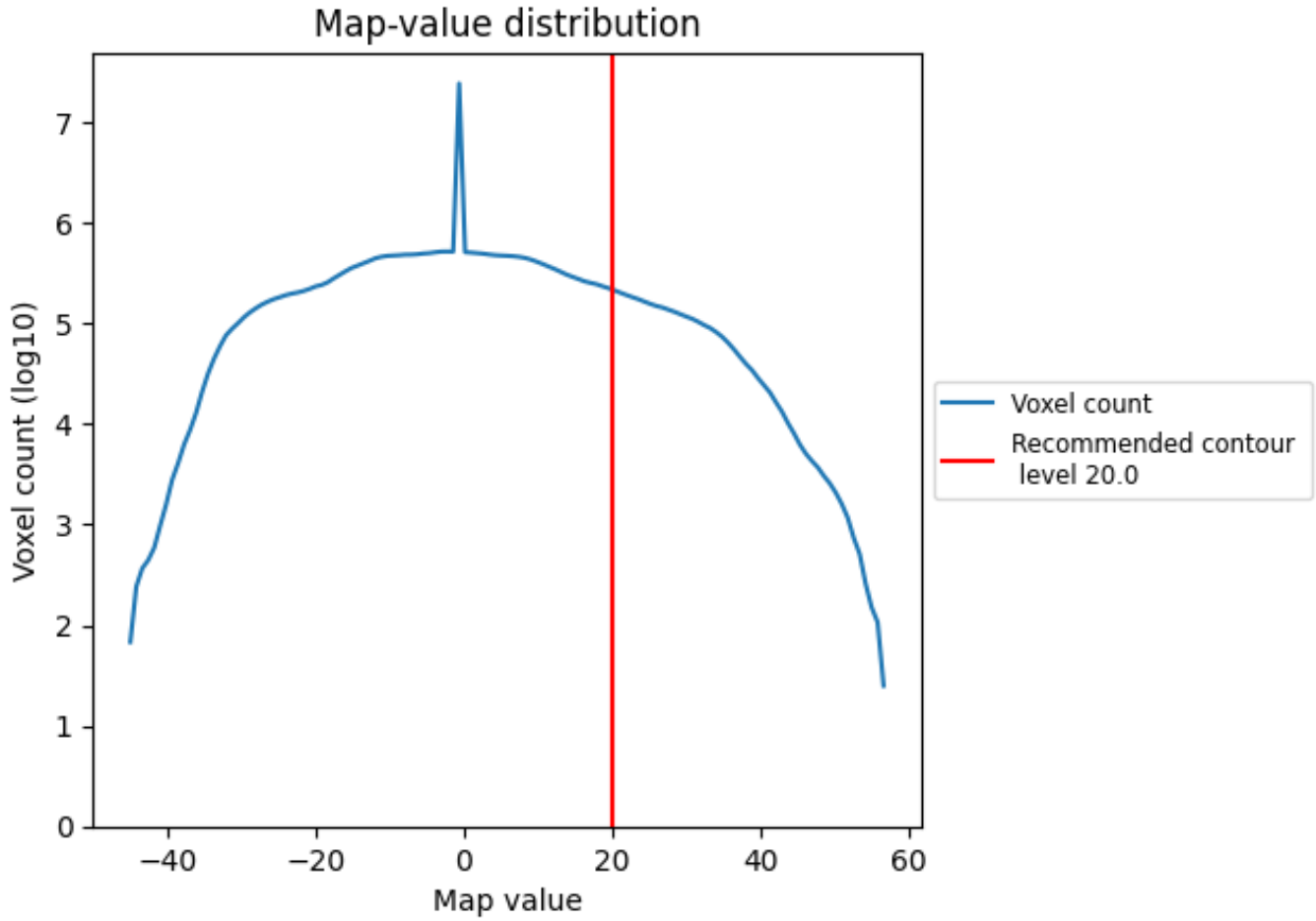
6.5 Mask visualisation (i)

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

7 Map analysis

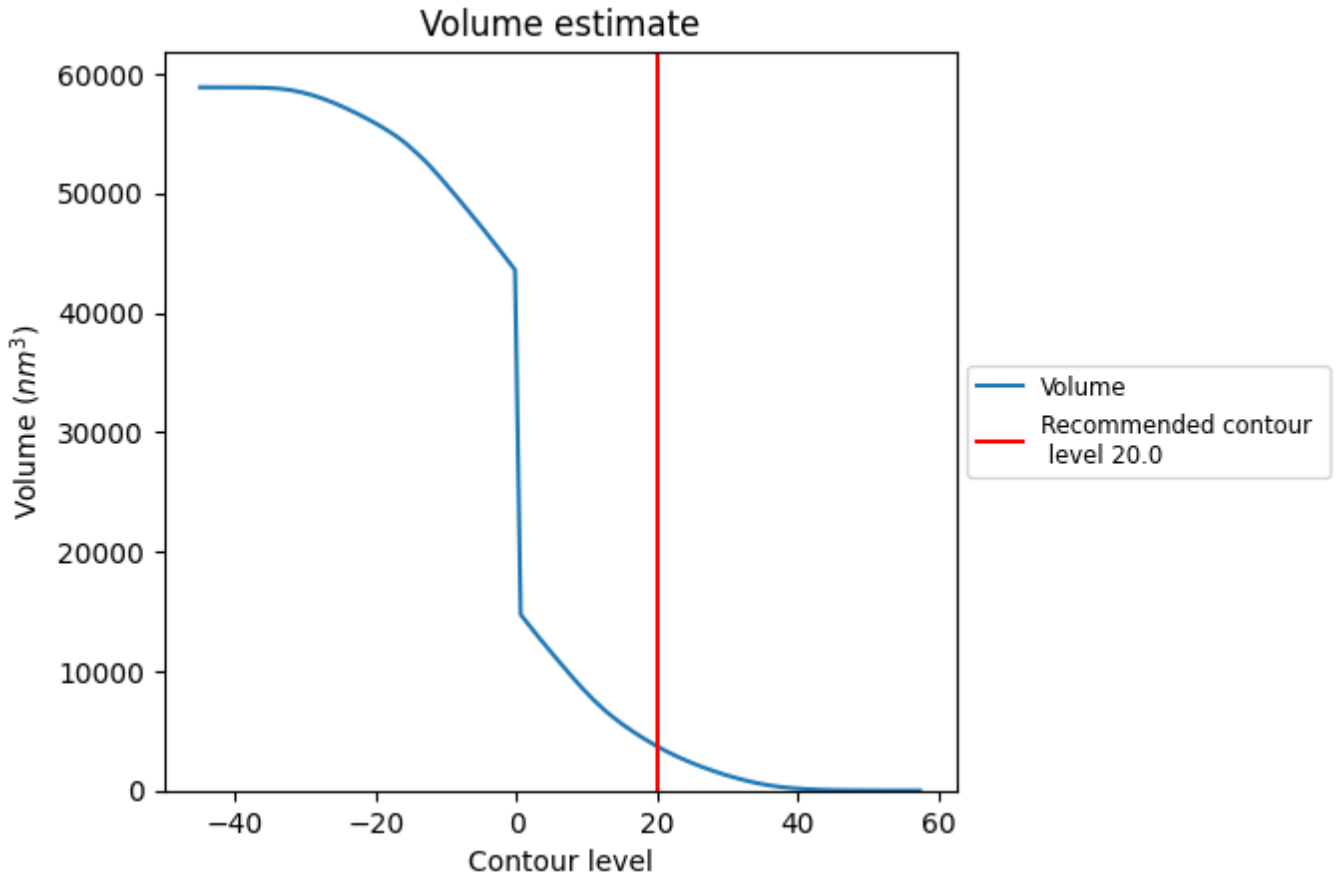
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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



The volume at the recommended contour level is 3713 nm^3; this corresponds to an approximate mass of 3354 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

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

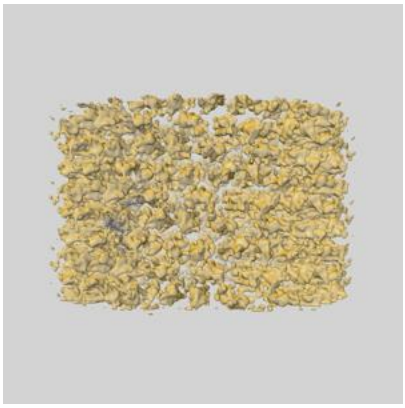
8 Fourier-Shell correlation (i)

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

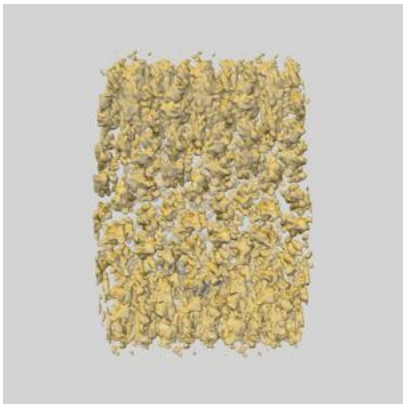
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-3076 and PDB model 5FJB. Per-residue inclusion information can be found in section 3 on page 4.

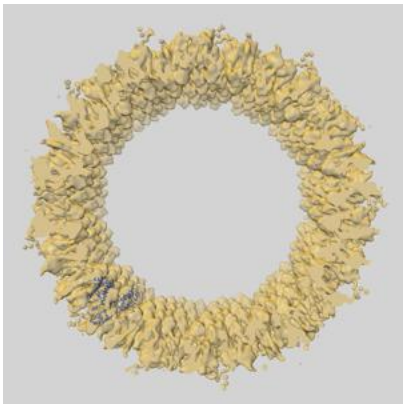
9.1 Map-model overlay



X



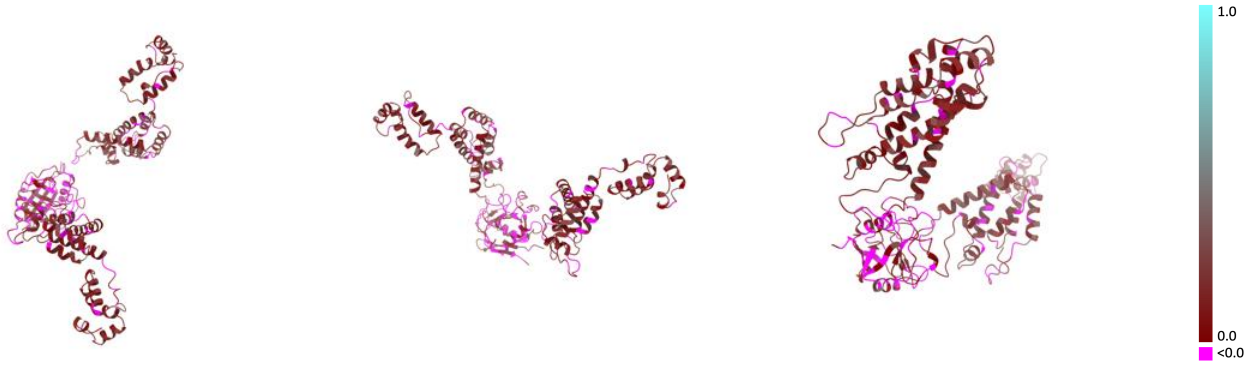
Y



Z

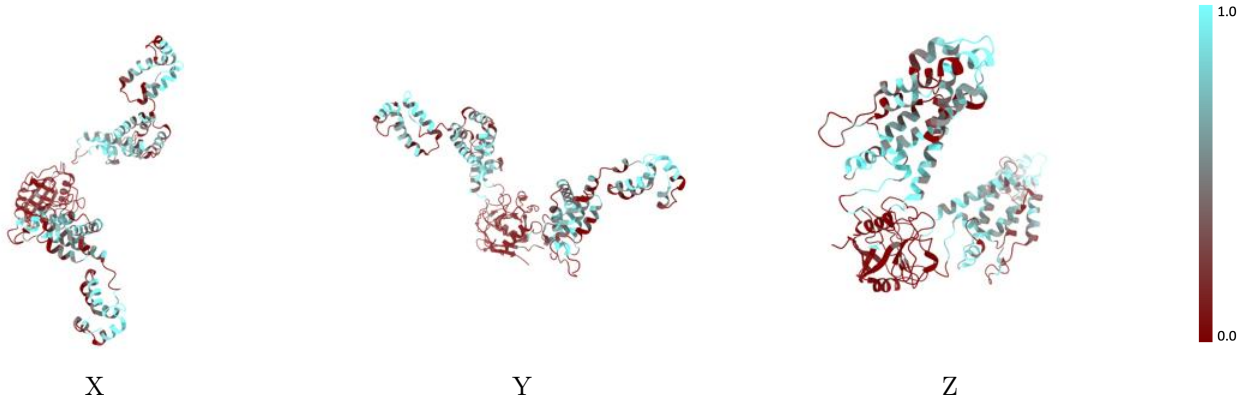
The images above show the 3D surface view of the map at the recommended contour level 20.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model (i)



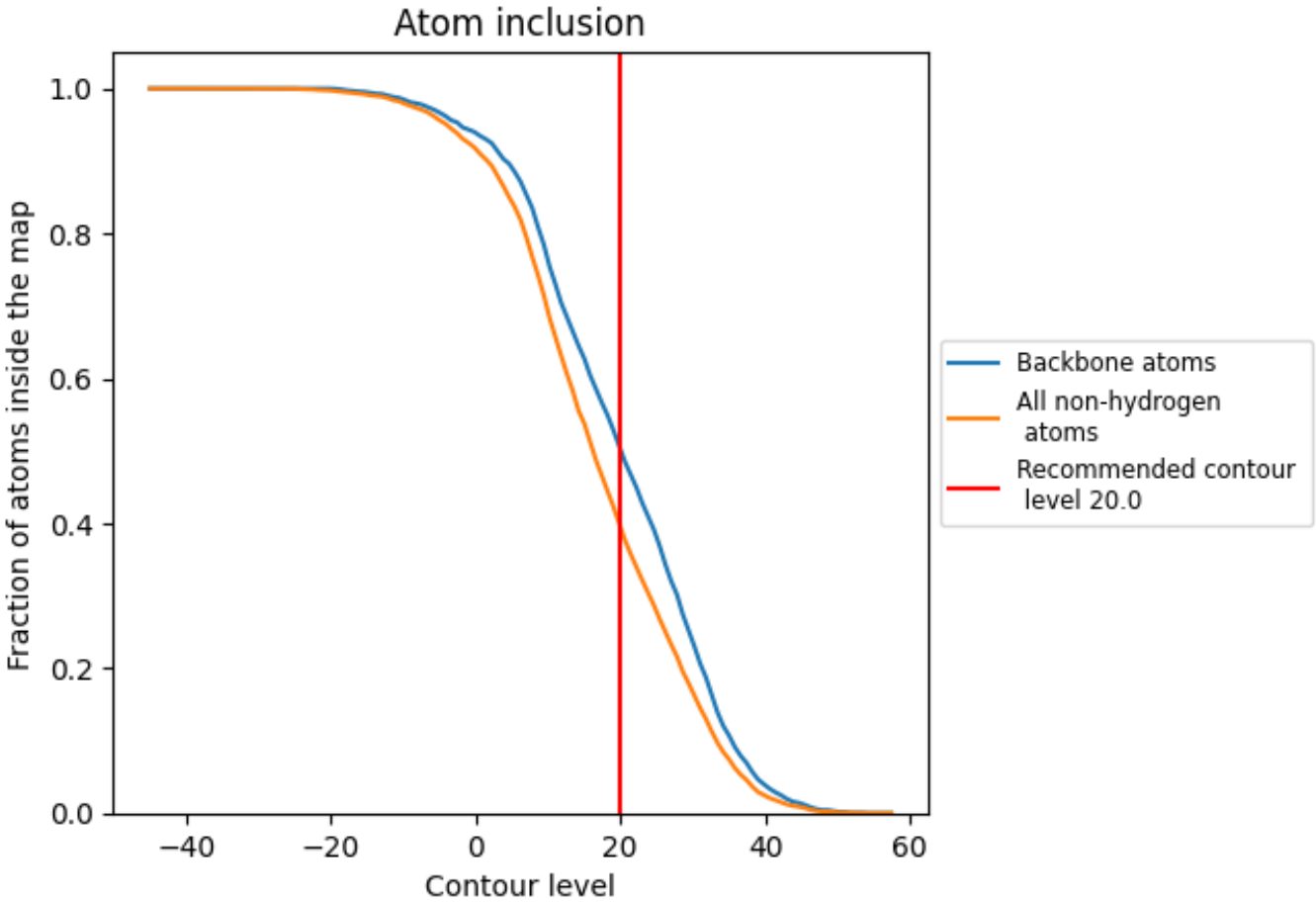
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (20.0).

9.4 Atom inclusion (i)



At the recommended contour level, 50% of all backbone atoms, 39% 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 (20.0) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.3945	0.1010
A	0.5269	0.1310
B	0.5352	0.1350
C	0.0265	0.0140