wwPDB EM Validation Summary Report

Oct 27, 2026 – 04:48 AM JST

PDB ID	:	44DL / pdb_000044dl
EMDB ID	:	EMD-82523
Title	:	Cryo-EM Structure of Tetrameric EcCaspase Filament in Apo State
Authors	:	Hao, F.; Xin, X.
Deposited on	:	2026-07-25
Resolution	:	3.01 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:

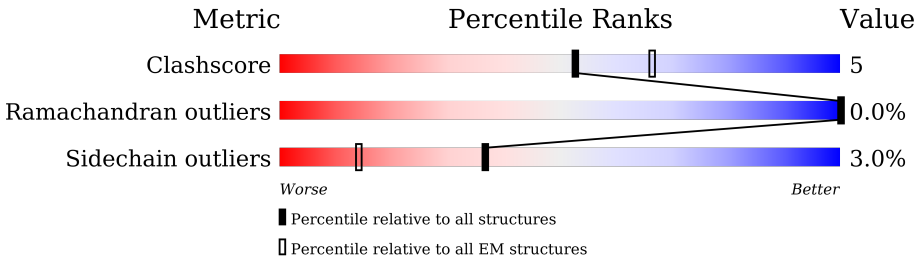
EMDB validation analysis	:	0.0.1.dev133
MolProbity	:	4-5-2 with Phenix2.2
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

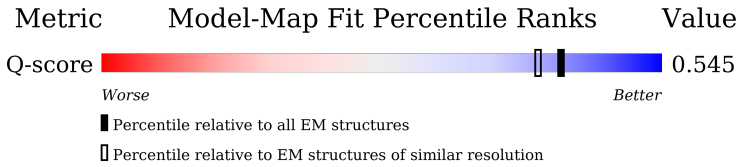
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.01 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.





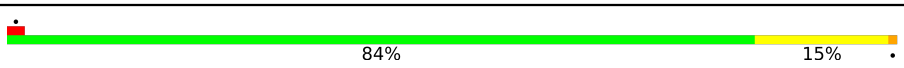
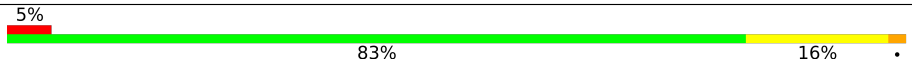
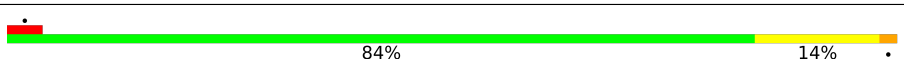
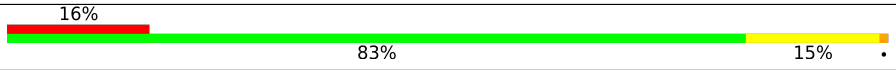
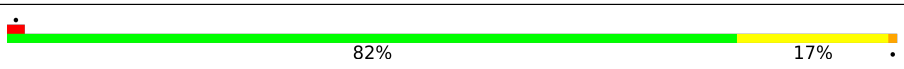
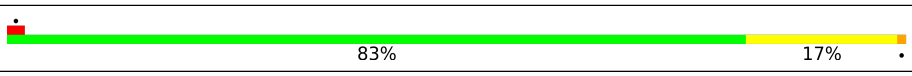
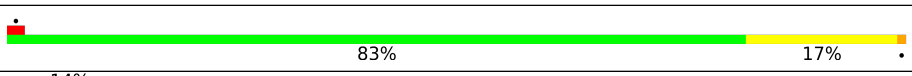
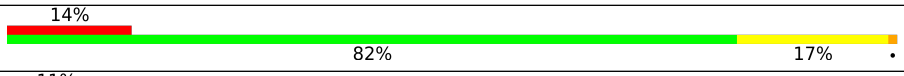
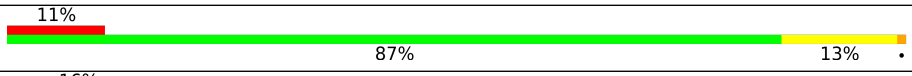
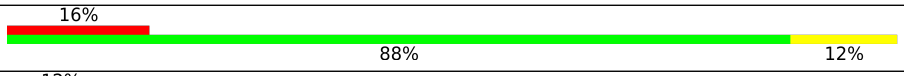
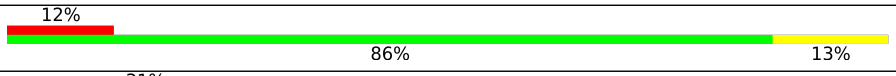
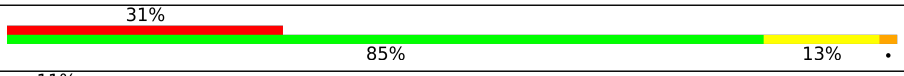
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13882 (2.51 - 3.51)

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	321	84% 15%
1	B	321	83% 16%
1	C	321	10% 88% 12%
1	D	321	11% 88% 11%

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Mol	Chain	Length	Quality of chain
1	E	321	
1	F	321	
1	G	321	
1	H	321	
1	I	321	
1	J	321	
1	K	321	
1	L	321	
1	M	321	
1	N	321	
1	O	321	
1	P	321	
1	Q	321	
1	R	321	
1	S	321	
1	T	321	

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 49980 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 Caspase family protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	A	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	D	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	C	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	I	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	E	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	Q	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	M	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	J	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	F	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	R	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	N	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	K	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	G	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	S	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	O	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	L	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	H	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	T	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		
1	P	321	Total	C	N	O	S	0	0
			2499	1579	432	473	15		

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	21	ALA	CYS	conflict	UNP A0AAF0KJ23
B	22	CYS	GLU	conflict	UNP A0AAF0KJ23
A	21	ALA	CYS	conflict	UNP A0AAF0KJ23
A	22	CYS	GLU	conflict	UNP A0AAF0KJ23
D	21	ALA	CYS	conflict	UNP A0AAF0KJ23
D	22	CYS	GLU	conflict	UNP A0AAF0KJ23
C	21	ALA	CYS	conflict	UNP A0AAF0KJ23
C	22	CYS	GLU	conflict	UNP A0AAF0KJ23
I	21	ALA	CYS	conflict	UNP A0AAF0KJ23
I	22	CYS	GLU	conflict	UNP A0AAF0KJ23
E	21	ALA	CYS	conflict	UNP A0AAF0KJ23
E	22	CYS	GLU	conflict	UNP A0AAF0KJ23
Q	21	ALA	CYS	conflict	UNP A0AAF0KJ23
Q	22	CYS	GLU	conflict	UNP A0AAF0KJ23
M	21	ALA	CYS	conflict	UNP A0AAF0KJ23
M	22	CYS	GLU	conflict	UNP A0AAF0KJ23
J	21	ALA	CYS	conflict	UNP A0AAF0KJ23
J	22	CYS	GLU	conflict	UNP A0AAF0KJ23
F	21	ALA	CYS	conflict	UNP A0AAF0KJ23
F	22	CYS	GLU	conflict	UNP A0AAF0KJ23
R	21	ALA	CYS	conflict	UNP A0AAF0KJ23
R	22	CYS	GLU	conflict	UNP A0AAF0KJ23
N	21	ALA	CYS	conflict	UNP A0AAF0KJ23
N	22	CYS	GLU	conflict	UNP A0AAF0KJ23
K	21	ALA	CYS	conflict	UNP A0AAF0KJ23
K	22	CYS	GLU	conflict	UNP A0AAF0KJ23
G	21	ALA	CYS	conflict	UNP A0AAF0KJ23
G	22	CYS	GLU	conflict	UNP A0AAF0KJ23
S	21	ALA	CYS	conflict	UNP A0AAF0KJ23
S	22	CYS	GLU	conflict	UNP A0AAF0KJ23
O	21	ALA	CYS	conflict	UNP A0AAF0KJ23
O	22	CYS	GLU	conflict	UNP A0AAF0KJ23

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Chain	Residue	Modelled	Actual	Comment	Reference
L	21	ALA	CYS	conflict	UNP A0AAF0KJ23
L	22	CYS	GLU	conflict	UNP A0AAF0KJ23
H	21	ALA	CYS	conflict	UNP A0AAF0KJ23
H	22	CYS	GLU	conflict	UNP A0AAF0KJ23
T	21	ALA	CYS	conflict	UNP A0AAF0KJ23
T	22	CYS	GLU	conflict	UNP A0AAF0KJ23
P	21	ALA	CYS	conflict	UNP A0AAF0KJ23
P	22	CYS	GLU	conflict	UNP A0AAF0KJ23

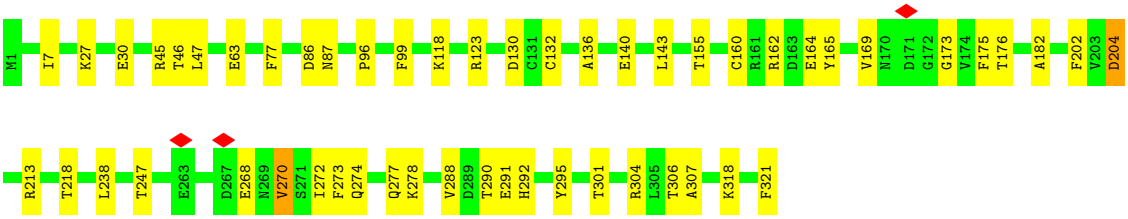
3 Residue-property plots (i)

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: Caspase family protein

Chain B:

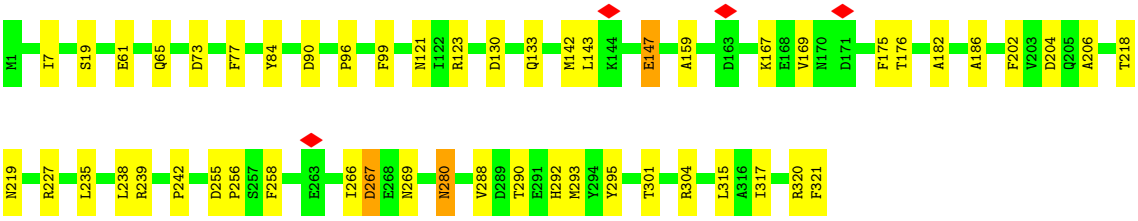




• Molecule 1: Caspase family protein

Chain A:

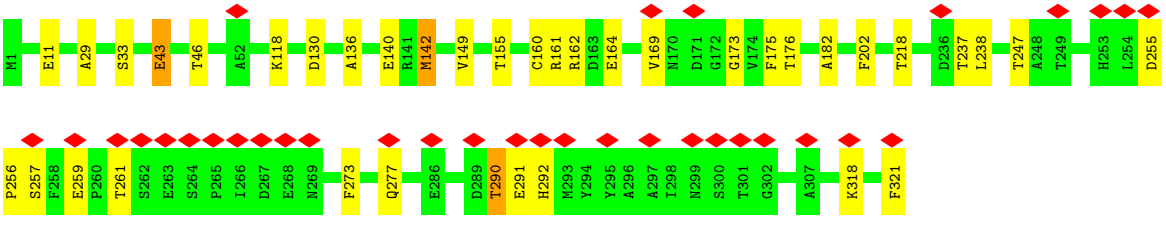




• Molecule 1: Caspase family protein

Chain D:

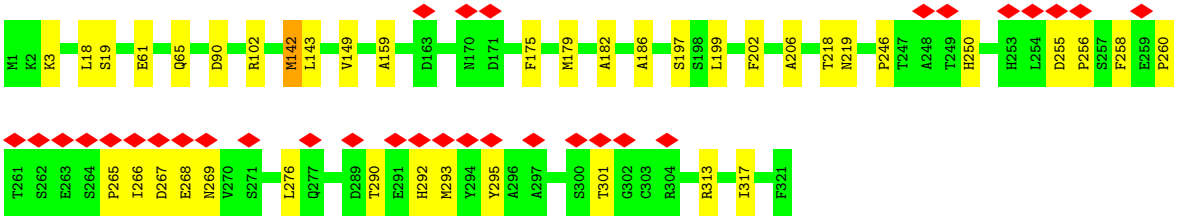




• Molecule 1: Caspase family protein

Chain C:

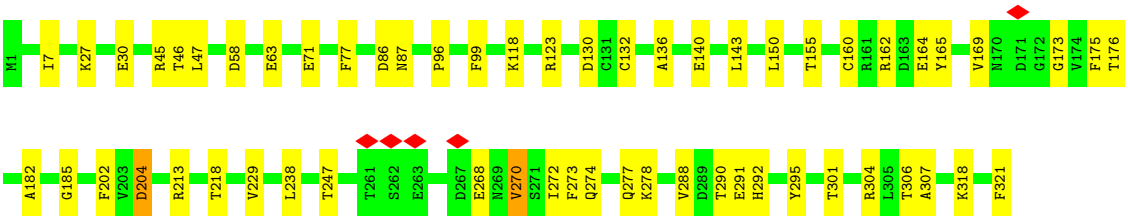




• Molecule 1: Caspase family protein

Chain I:

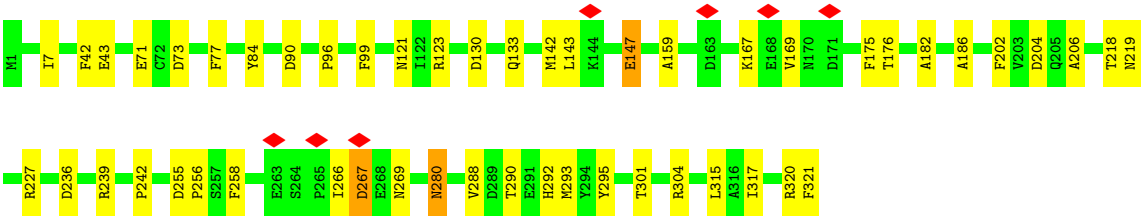




• Molecule 1: Caspase family protein

Chain E:

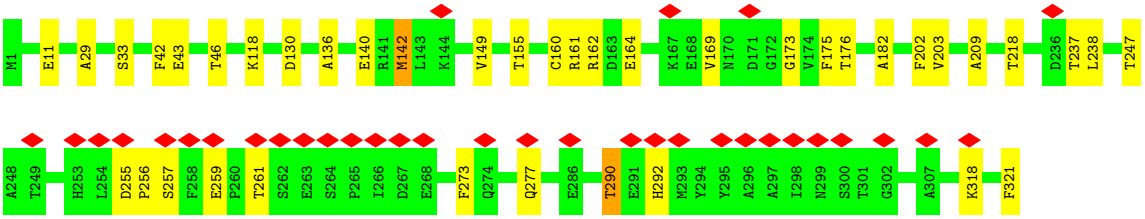




• Molecule 1: Caspase family protein

Chain Q:

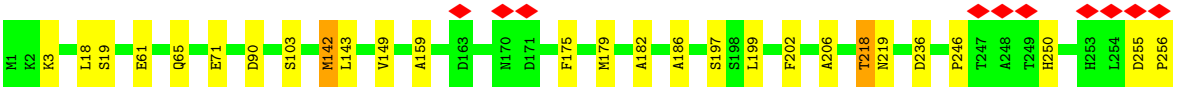


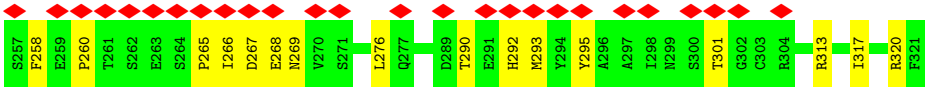


• Molecule 1: Caspase family protein

Chain M:



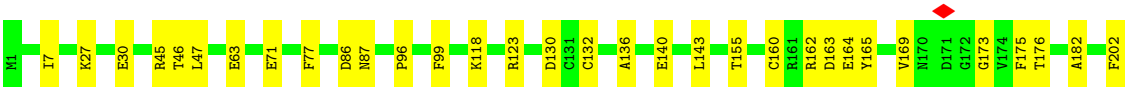




• Molecule 1: Caspase family protein

Chain J:





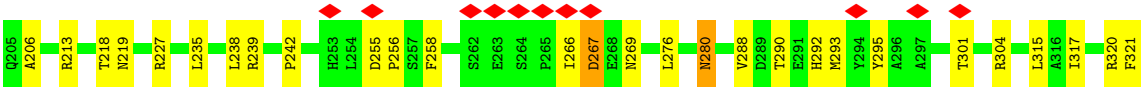


• Molecule 1: Caspase family protein

Chain F:



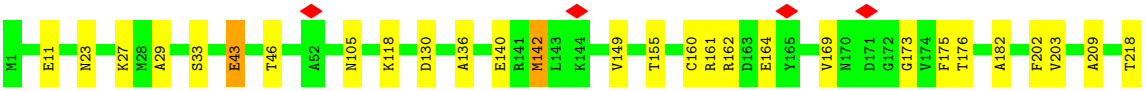


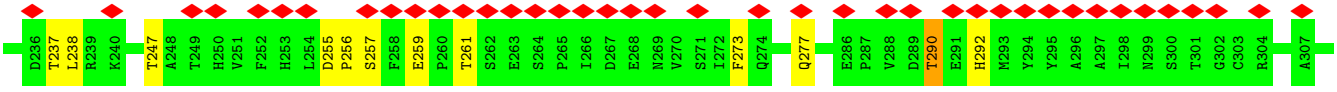


• Molecule 1: Caspase family protein

Chain R:







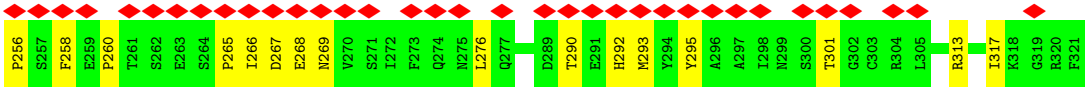


• Molecule 1: Caspase family protein

Chain N:



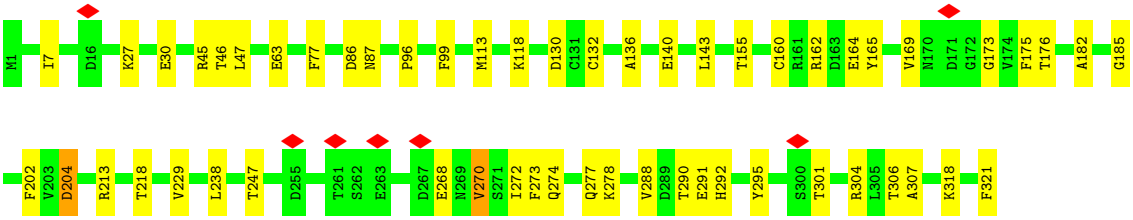




• Molecule 1: Caspase family protein

Chain K:

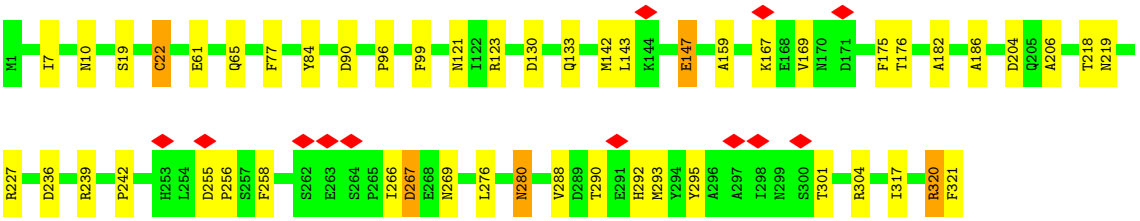




• Molecule 1: Caspase family protein

Chain G:

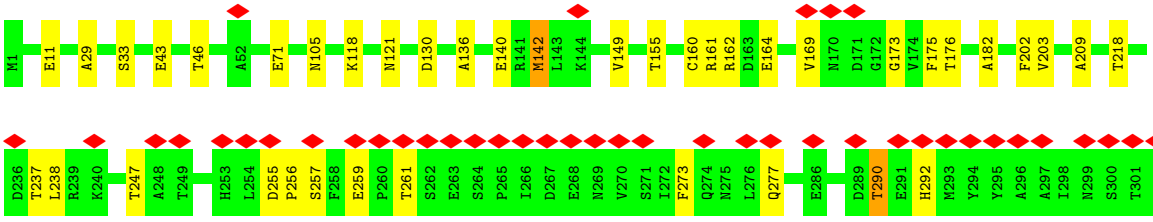




• Molecule 1: Caspase family protein

Chain S:

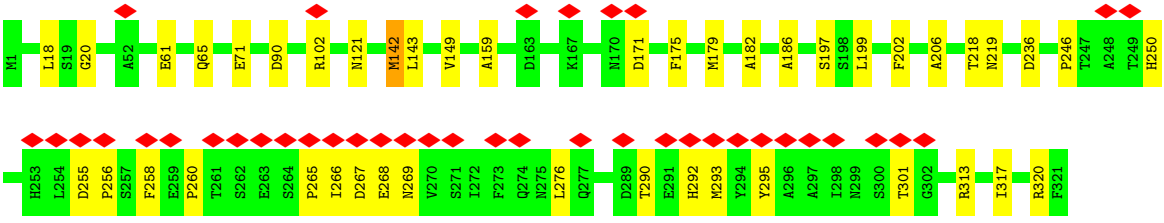




• Molecule 1: Caspase family protein

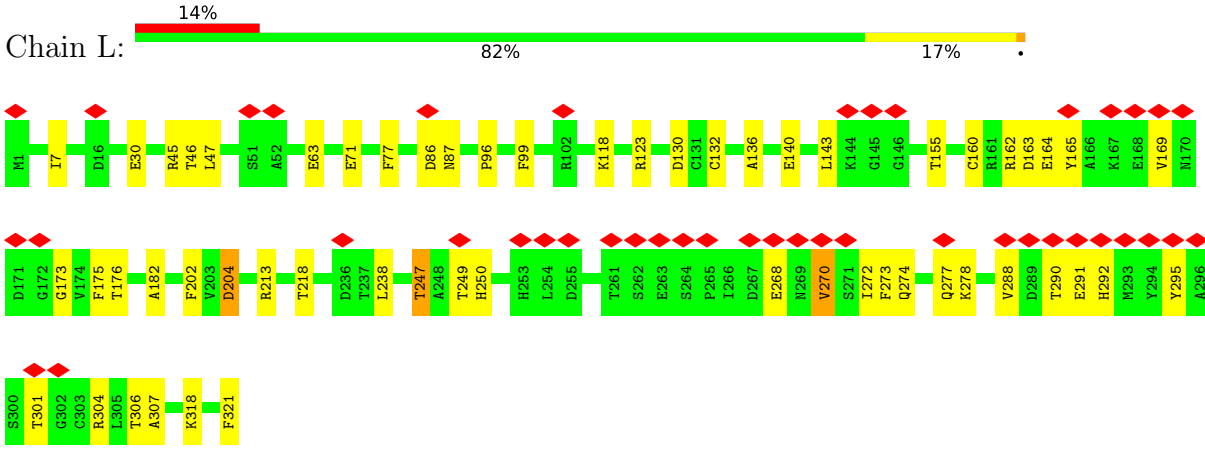
Chain O:





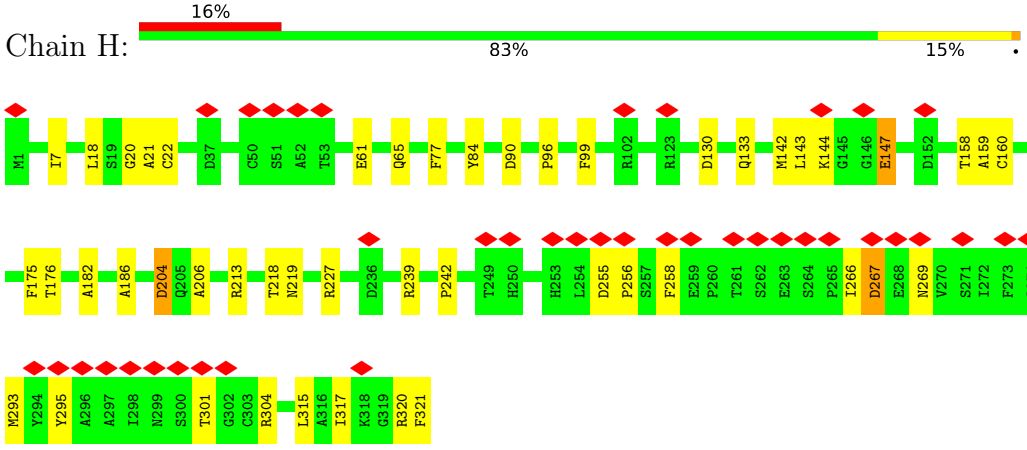
• Molecule 1: Caspase family protein

Chain L:



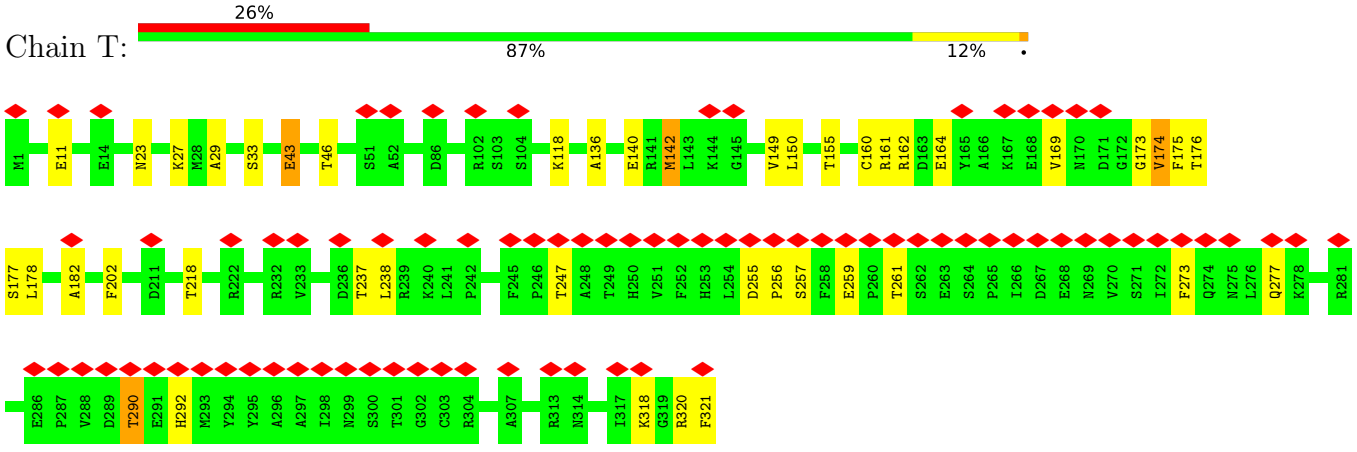
• Molecule 1: Caspase family protein

Chain H:



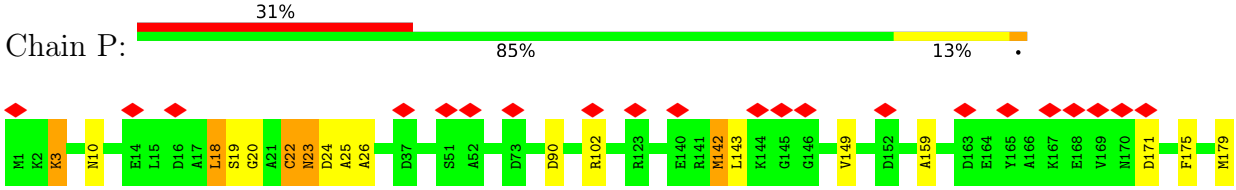
• Molecule 1: Caspase family protein

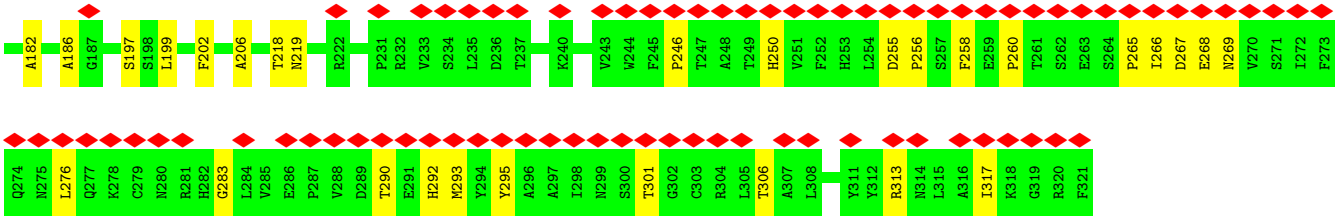
Chain T:



• Molecule 1: Caspase family protein

Chain P:




A182
A186
G187
S197
S198
L199
F202
A206
T218
N219
R222
P231
R232
V233
S234
L235
D236
T237
K240
V243
W244
F245
P246
T247
A248
T249
H250
V251
F252
H253
L254
D255
P256
S257
F258
E259
P260
T261
S262
E263
S264
P265
I266
D267
E268
N269
V270
S271
I272
F273
Q274
N275
L276
Q277
K278
C279
N280
R281
H282
G283
L284
V285
E286
P287
V288
D289
T290
E291
H292
M293
Y294
Y295
A296
A297
I298
N299
S300
T301
G302
C303
R304
L305
T306
A307
L308
Y311
Y312
R313
N314
L315
A316
I317
K318
G319
R320
F321

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	20142	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	45	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.666	Depositor
Minimum map value	-0.403	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.026	Depositor
Recommended contour level	0.09	Depositor
Map size (Å)	317.4144, 317.4144, 317.4144	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8266, 0.8266, 0.8266	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	0.17	0/2554	0.33	0/3462
1	B	0.14	0/2554	0.28	0/3462
1	C	0.14	0/2554	0.31	0/3462
1	D	0.14	0/2554	0.30	0/3462
1	E	0.17	0/2554	0.35	1/3462 (0.0%)
1	F	0.16	0/2554	0.32	0/3462
1	G	0.26	0/2554	0.39	1/3462 (0.0%)
1	H	0.20	0/2554	0.34	1/3462 (0.0%)
1	I	0.15	0/2554	0.29	0/3462
1	J	0.15	0/2554	0.29	0/3462
1	K	0.15	0/2554	0.28	0/3462
1	L	0.15	0/2554	0.29	0/3462
1	M	0.16	0/2554	0.33	0/3462
1	N	0.14	0/2554	0.31	0/3462
1	O	0.24	0/2554	0.36	1/3462 (0.0%)
1	P	0.22	0/2554	0.36	2/3462 (0.1%)
1	Q	0.17	0/2554	0.33	1/3462 (0.0%)
1	R	0.15	0/2554	0.31	0/3462
1	S	0.13	0/2554	0.29	0/3462
1	T	0.21	0/2554	0.32	0/3462
All	All	0.17	0/51080	0.32	7/69240 (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	G	0	2
1	M	0	1
1	O	0	1
1	R	0	1
1	T	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
All	All	0	6

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	42	PHE	CA-CB-CG	7.32	121.12	113.80
1	G	22	CYS	N-CA-C	-7.05	106.59	114.62
1	O	18	LEU	N-CA-C	-6.48	100.38	110.17
1	P	23	ASN	N-CA-C	-6.35	103.99	111.03
1	Q	42	PHE	CA-CB-CG	5.94	119.74	113.80

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	G	123	ARG	Sidechain
1	G	320	ARG	Sidechain
1	M	320	ARG	Sidechain
1	O	320	ARG	Sidechain
1	R	320	ARG	Sidechain

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	2499	0	2436	33	0
1	B	2499	0	2436	33	0
1	C	2499	0	2436	27	0
1	D	2499	0	2436	26	0
1	E	2499	0	2436	30	0
1	F	2499	0	2436	36	0
1	G	2499	0	2436	31	0
1	H	2499	0	2436	32	0
1	I	2499	0	2436	37	0
1	J	2499	0	2436	34	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	K	2499	0	2436	34	0
1	L	2499	0	2436	35	0
1	M	2499	0	2436	27	0
1	N	2499	0	2436	24	0
1	O	2499	0	2436	31	0
1	P	2499	0	2436	33	0
1	Q	2499	0	2436	26	0
1	R	2499	0	2436	32	0
1	S	2499	0	2436	28	0
1	T	2499	0	2436	29	0
All	All	49980	0	48720	520	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 5.

The worst 5 of 520 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:D:43:GLU:OE2	1:S:209:ALA:CB	1.85	1.23
1:Q:209:ALA:CB	1:R:43:GLU:OE2	1.86	1.21
1:R:209:ALA:CB	1:T:43:GLU:OE2	1.89	1.20
1:O:71:GLU:HG2	1:O:121:ASN:HB3	1.20	1.16
1:Q:209:ALA:HB3	1:R:43:GLU:OE2	1.44	1.15

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	319/321 (99%)	302 (95%)	17 (5%)	0	100	100
1	B	319/321 (99%)	308 (97%)	11 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	319/321 (99%)	300 (94%)	19 (6%)	0	100	100
1	D	319/321 (99%)	308 (97%)	11 (3%)	0	100	100
1	E	319/321 (99%)	301 (94%)	17 (5%)	1 (0%)	36	68
1	F	319/321 (99%)	302 (95%)	17 (5%)	0	100	100
1	G	319/321 (99%)	302 (95%)	17 (5%)	0	100	100
1	H	319/321 (99%)	303 (95%)	16 (5%)	0	100	100
1	I	319/321 (99%)	308 (97%)	11 (3%)	0	100	100
1	J	319/321 (99%)	308 (97%)	11 (3%)	0	100	100
1	K	319/321 (99%)	308 (97%)	11 (3%)	0	100	100
1	L	319/321 (99%)	307 (96%)	12 (4%)	0	100	100
1	M	319/321 (99%)	299 (94%)	20 (6%)	0	100	100
1	N	319/321 (99%)	301 (94%)	18 (6%)	0	100	100
1	O	319/321 (99%)	301 (94%)	18 (6%)	0	100	100
1	P	319/321 (99%)	300 (94%)	19 (6%)	0	100	100
1	Q	319/321 (99%)	308 (97%)	11 (3%)	0	100	100
1	R	319/321 (99%)	308 (97%)	11 (3%)	0	100	100
1	S	319/321 (99%)	308 (97%)	11 (3%)	0	100	100
1	T	319/321 (99%)	306 (96%)	13 (4%)	0	100	100
All	All	6380/6420 (99%)	6088 (95%)	291 (5%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	43	GLU

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	266/266 (100%)	259 (97%)	7 (3%)	40	71
1	B	266/266 (100%)	259 (97%)	7 (3%)	40	71
1	C	266/266 (100%)	260 (98%)	6 (2%)	44	72
1	D	266/266 (100%)	257 (97%)	9 (3%)	32	64
1	E	266/266 (100%)	257 (97%)	9 (3%)	32	64
1	F	266/266 (100%)	258 (97%)	8 (3%)	36	68
1	G	266/266 (100%)	257 (97%)	9 (3%)	32	64
1	H	266/266 (100%)	259 (97%)	7 (3%)	40	71
1	I	266/266 (100%)	259 (97%)	7 (3%)	40	71
1	J	266/266 (100%)	258 (97%)	8 (3%)	36	68
1	K	266/266 (100%)	259 (97%)	7 (3%)	40	71
1	L	266/266 (100%)	258 (97%)	8 (3%)	36	68
1	M	266/266 (100%)	257 (97%)	9 (3%)	32	64
1	N	266/266 (100%)	259 (97%)	7 (3%)	40	71
1	O	266/266 (100%)	261 (98%)	5 (2%)	50	75
1	P	266/266 (100%)	259 (97%)	7 (3%)	40	71
1	Q	266/266 (100%)	257 (97%)	9 (3%)	32	64
1	R	266/266 (100%)	257 (97%)	9 (3%)	32	64
1	S	266/266 (100%)	257 (97%)	9 (3%)	32	64
1	T	266/266 (100%)	256 (96%)	10 (4%)	29	62
All	All	5320/5320 (100%)	5163 (97%)	157 (3%)	37	68

5 of 157 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	S	155	THR
1	T	142	MET
1	S	290	THR
1	L	249	THR
1	P	3	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
1	L	250	HIS

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Mol	Chain	Res	Type
1	L	277	GLN
1	P	65	GLN
1	J	124	ASN
1	Q	124	ASN

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 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.

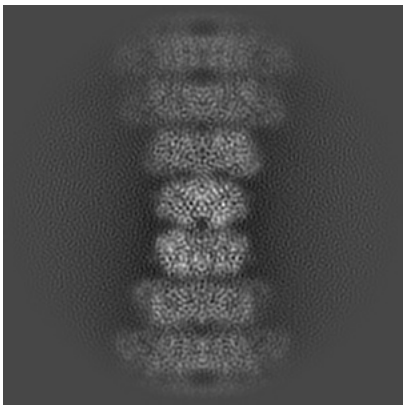
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-82523. These allow visual inspection of the internal detail of the map and identification of artifacts.

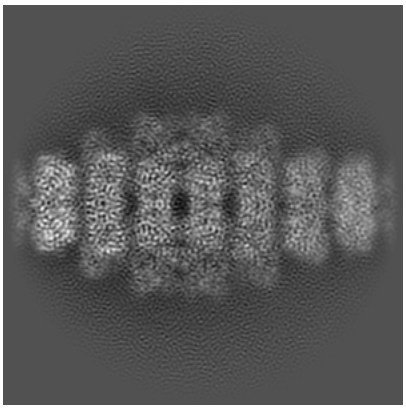
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

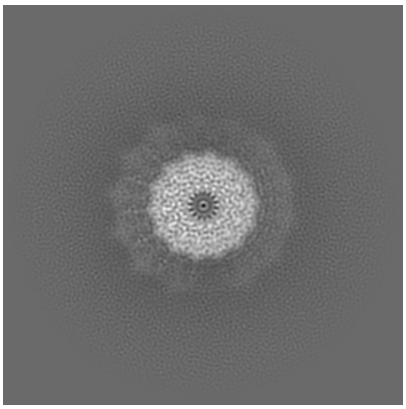
6.1.1 Primary map



X

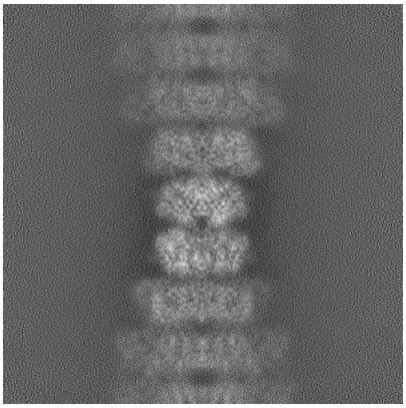


Y

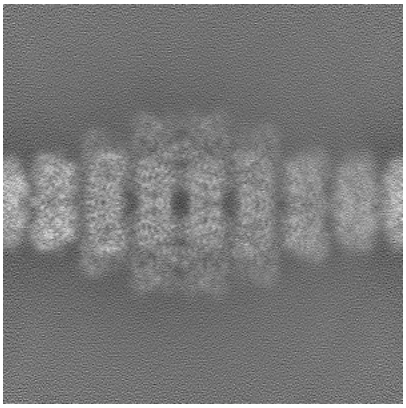


Z

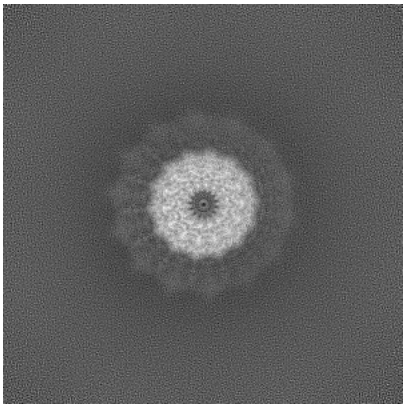
6.1.2 Raw map



X



Y

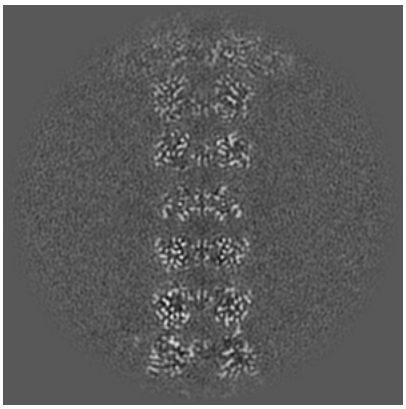


Z

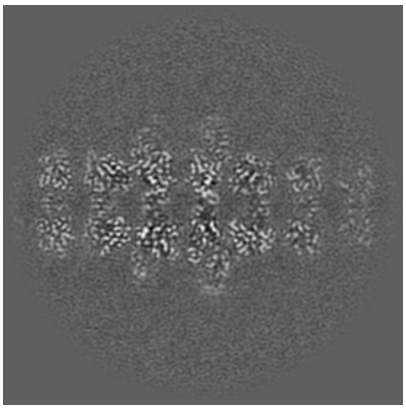
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

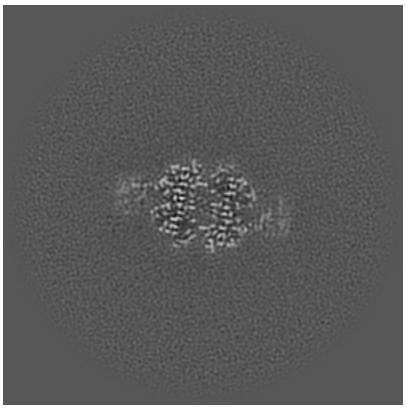
6.2.1 Primary map



X Index: 192

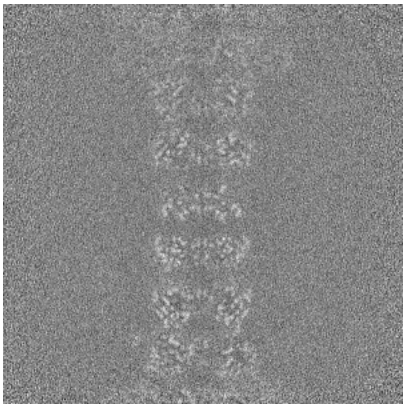


Y Index: 192

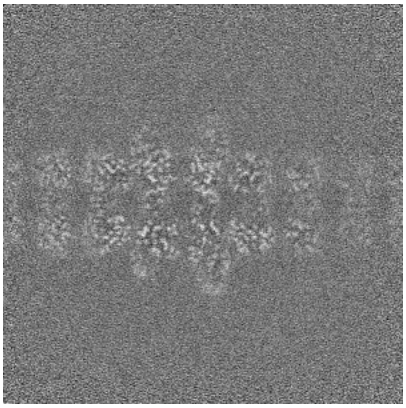


Z Index: 192

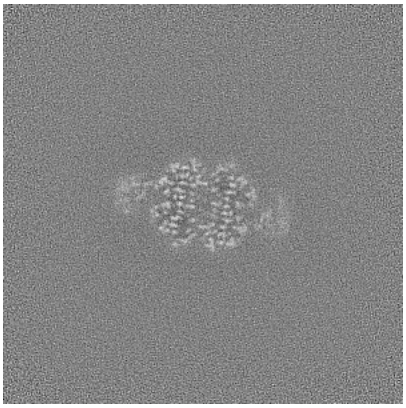
6.2.2 Raw map



X Index: 192



Y Index: 192

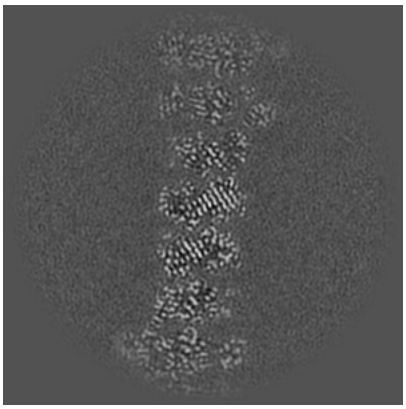


Z Index: 192

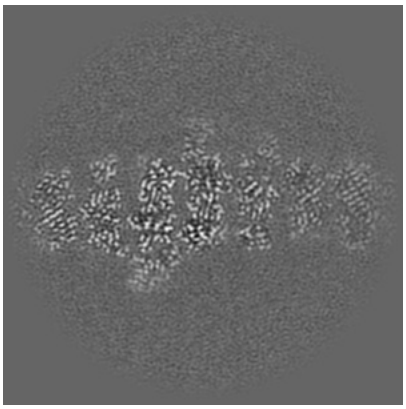
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

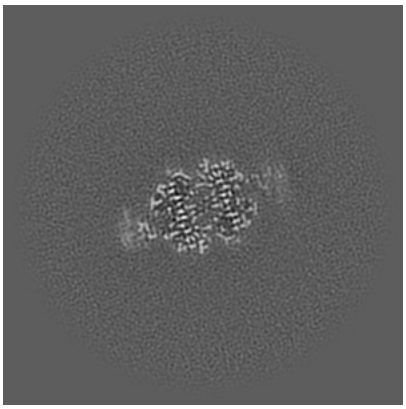
6.3.1 Primary map



X Index: 171

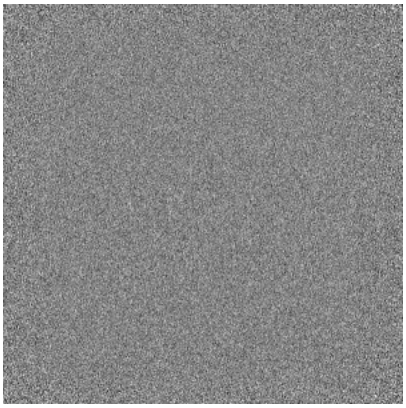


Y Index: 172

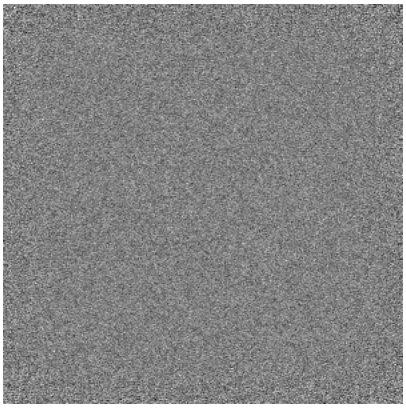


Z Index: 148

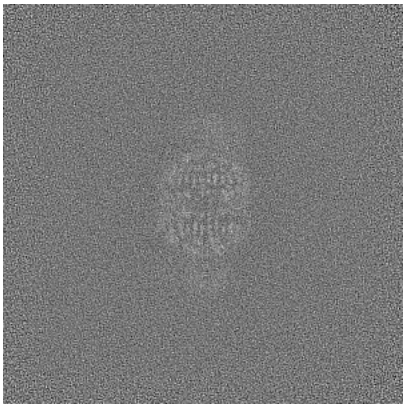
6.3.2 Raw map



X Index: 0



Y Index: 0

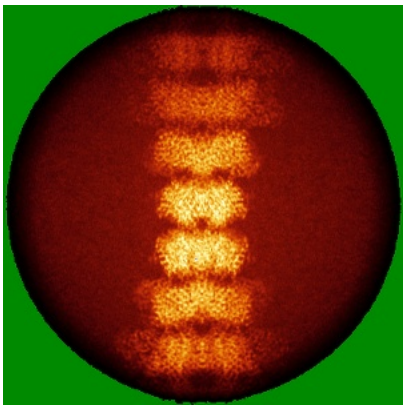


Z Index: 0

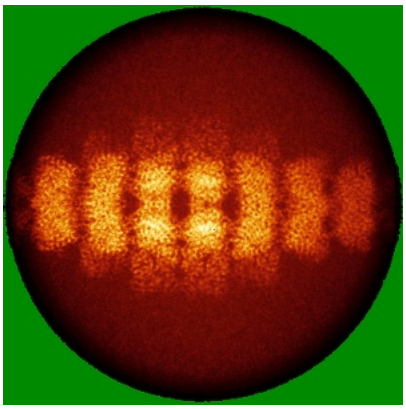
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

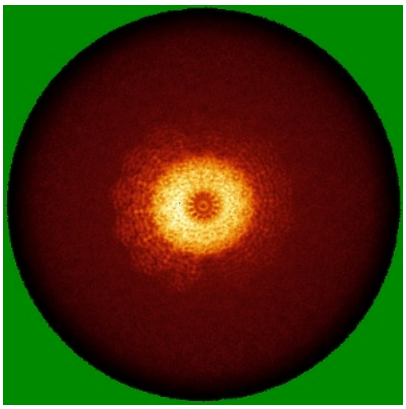
6.4.1 Primary map



X

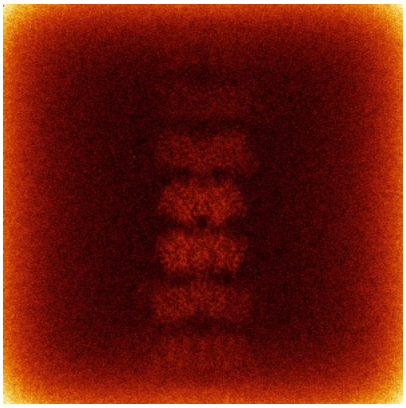


Y

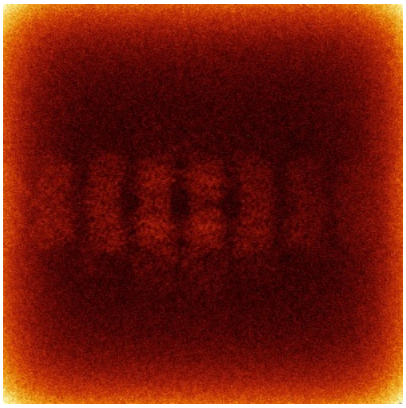


Z

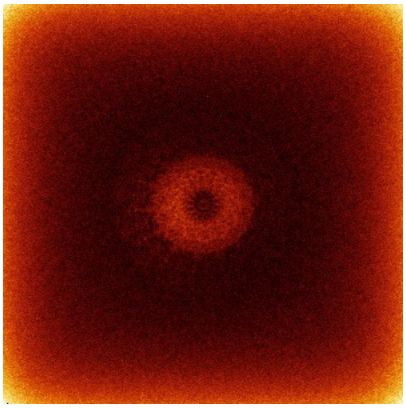
6.4.2 Raw map



X



Y

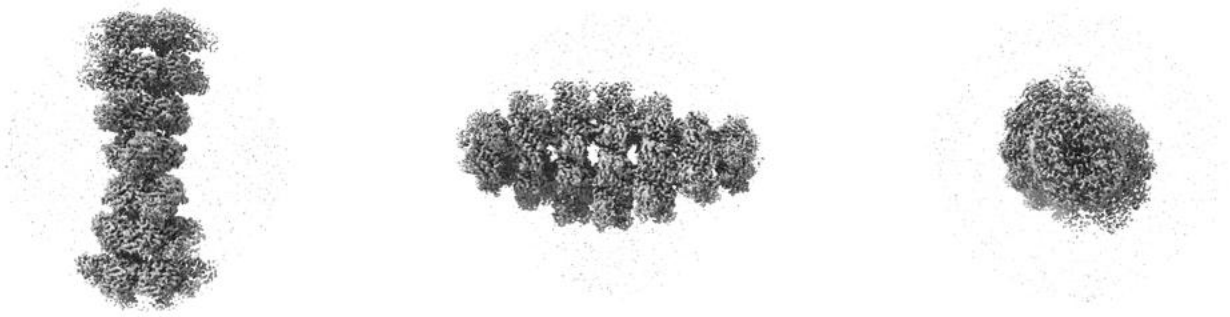


Z

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)

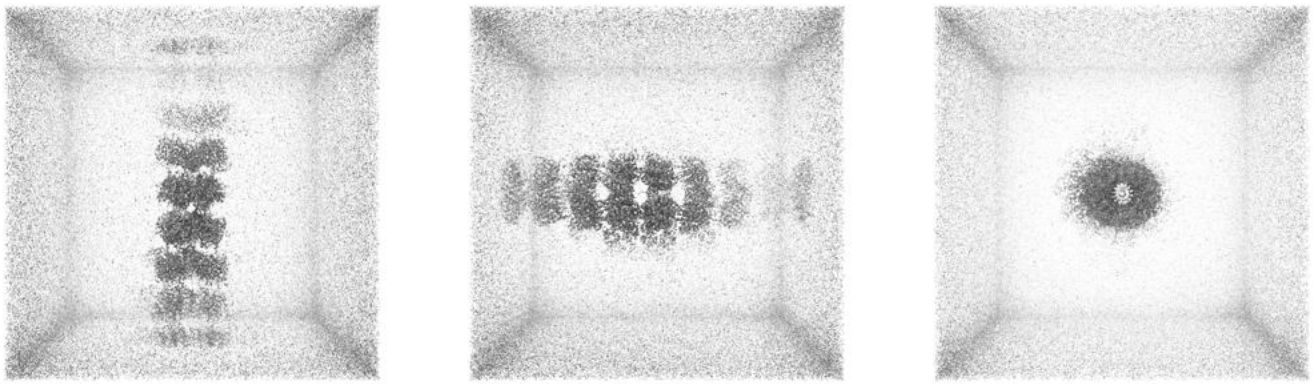
6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.09. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X Y Z

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.

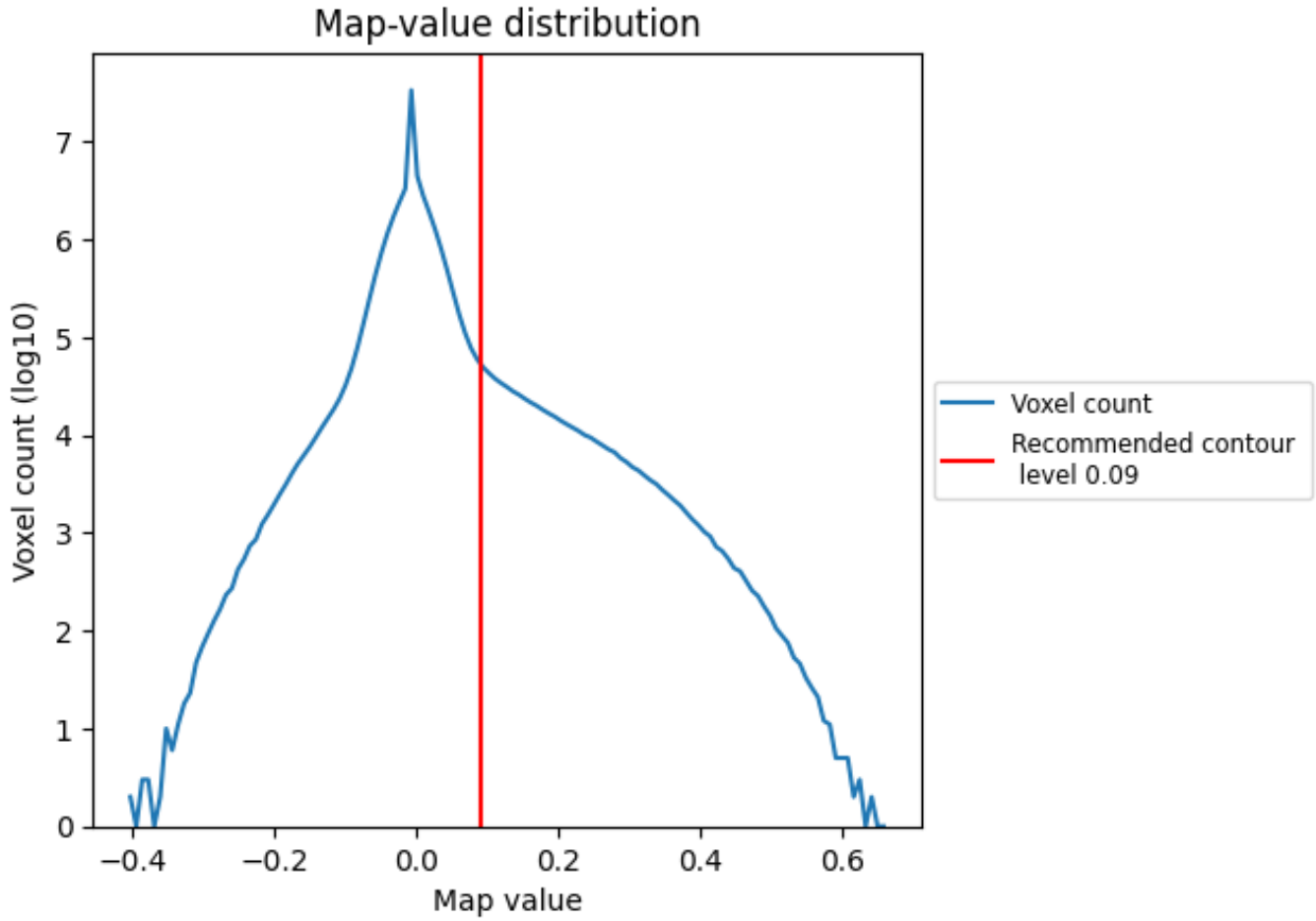
6.6 Mask visualisation (i)

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

7 Map analysis

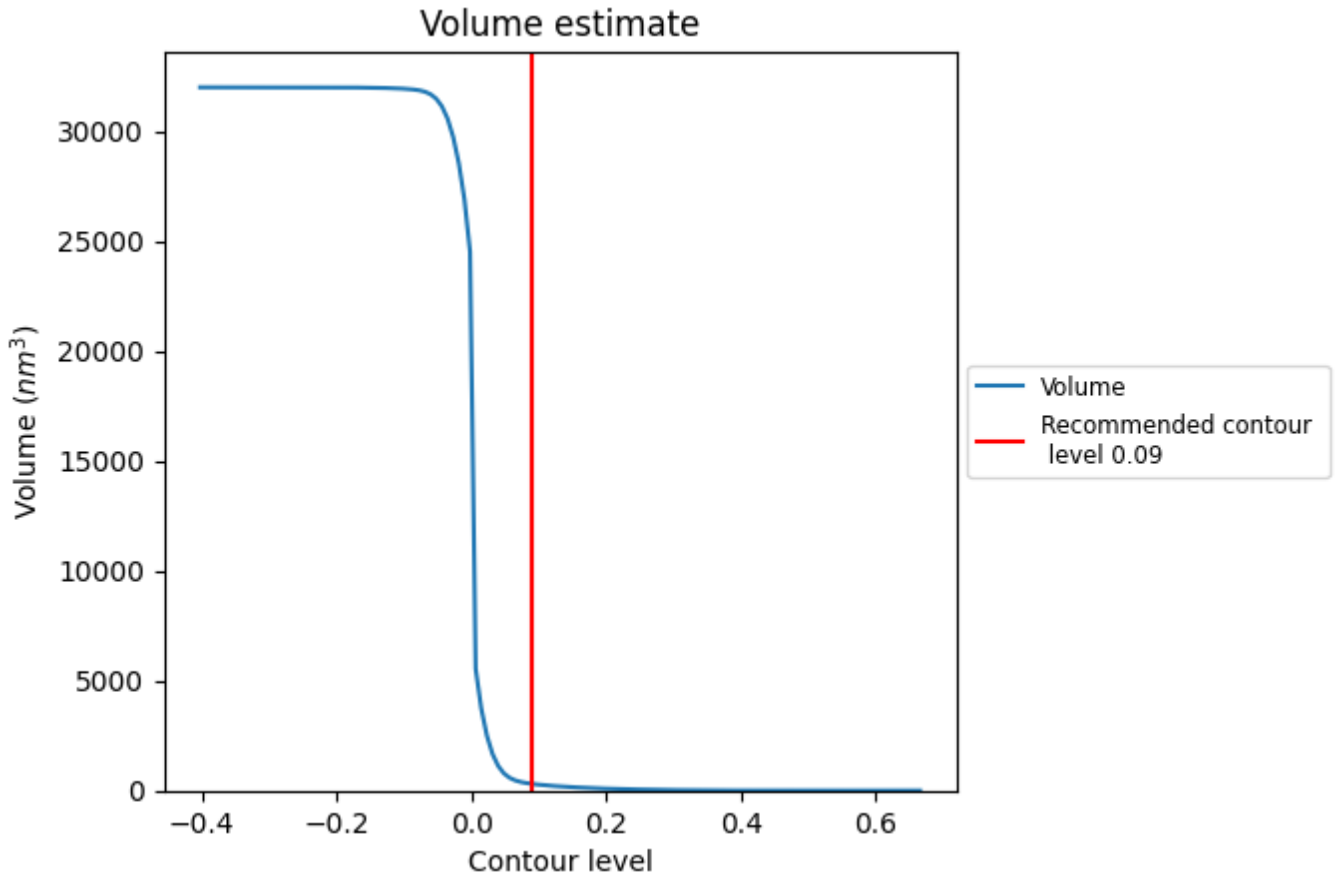
This section contains the results of statistical analysis of the map.

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.

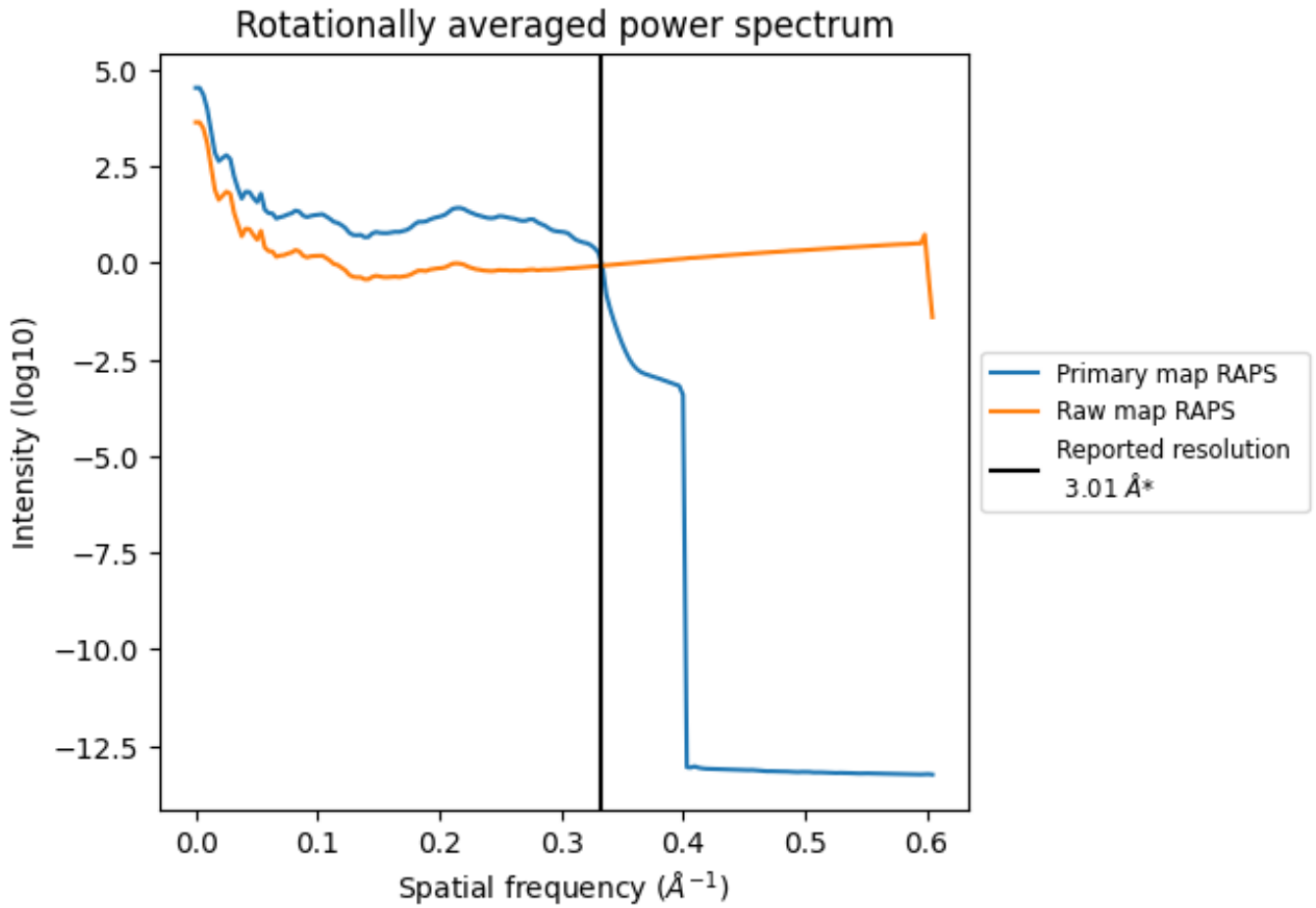
7.2 Volume estimate (i)



The volume at the recommended contour level is 303 nm^3; this corresponds to an approximate mass of 274 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)

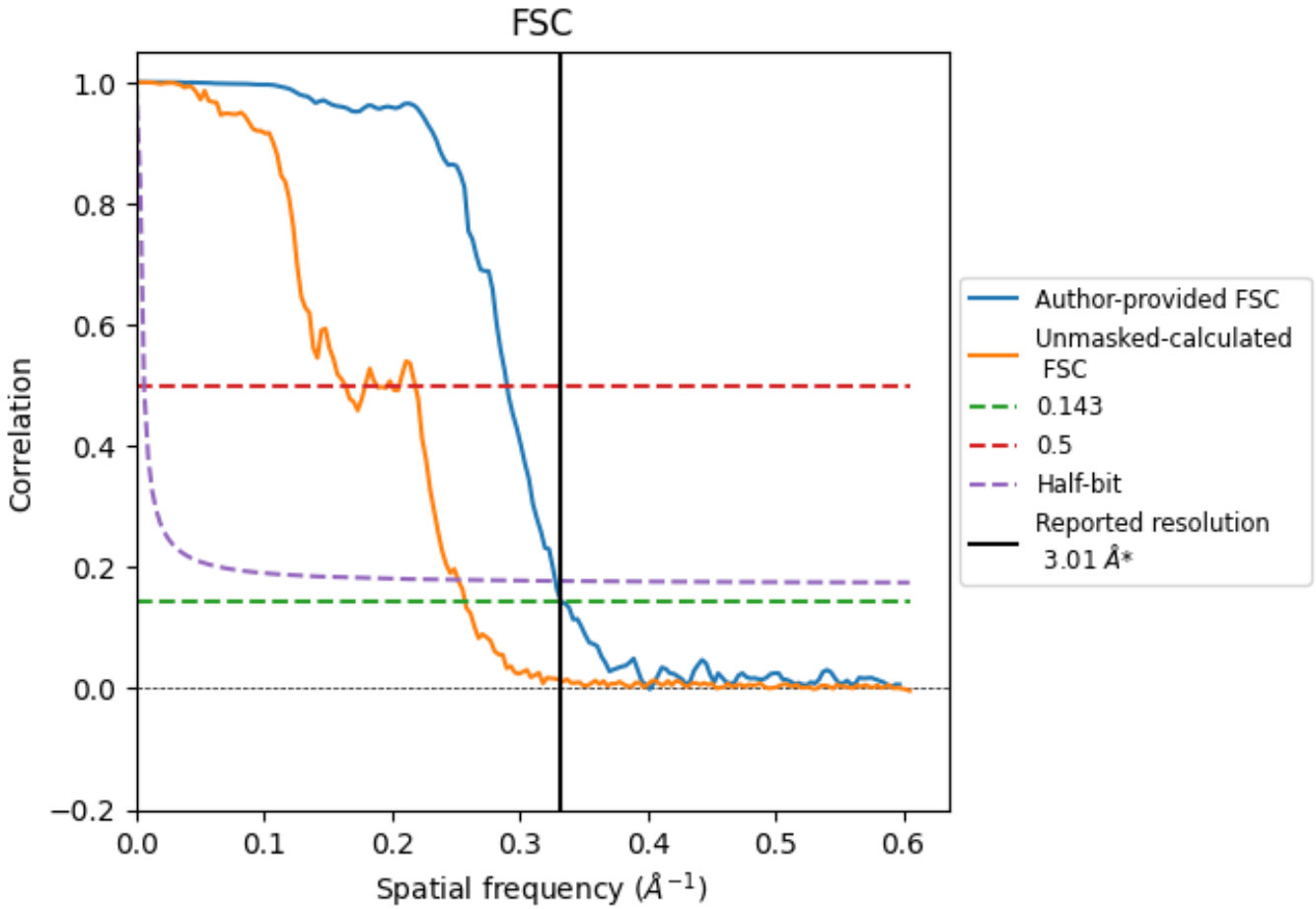


*Reported resolution corresponds to spatial frequency of 0.332 $Å^{-1}$

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.332 $Å^{-1}$

8.2 Resolution estimates (i)

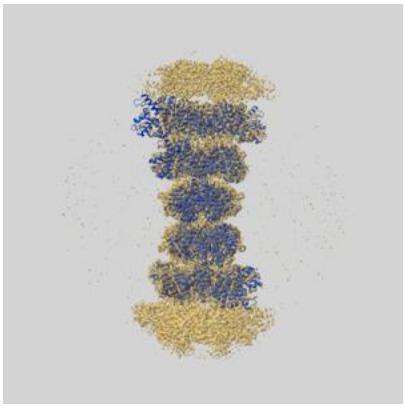
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.01	-	-
Author-provided FSC curve	3.01	3.45	3.05
Unmasked-calculated*	3.89	6.11	3.97

*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 3.89 differs from the reported value 3.01 by more than 10 %

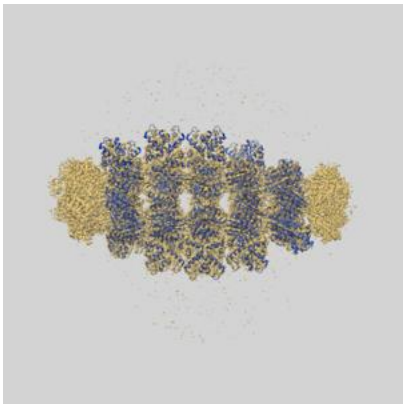
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-82523 and PDB model 44DL. Per-residue inclusion information can be found in section 3 on page 7.

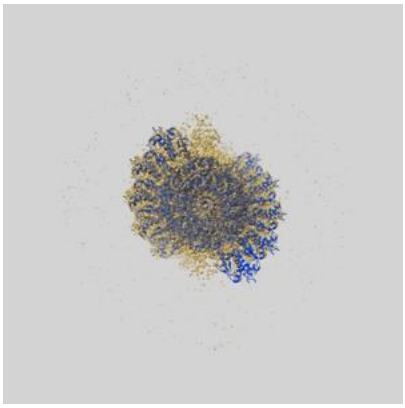
9.1 Map-model overlay



X



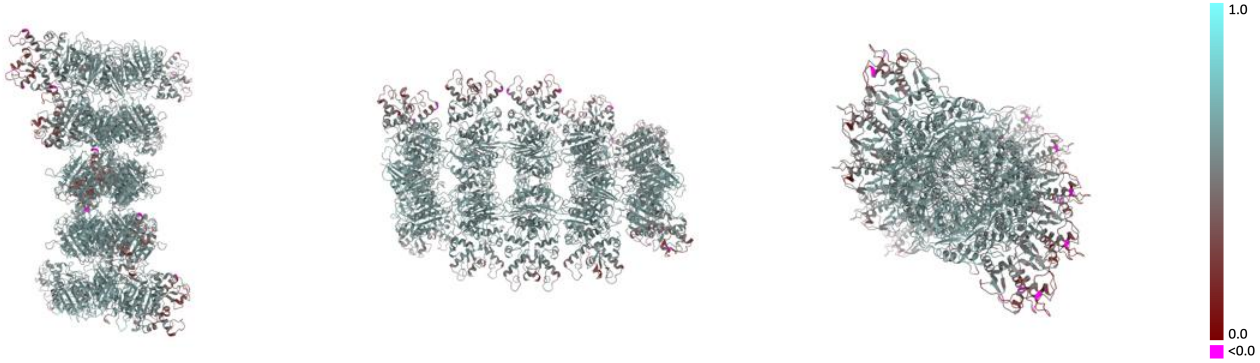
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.09 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.

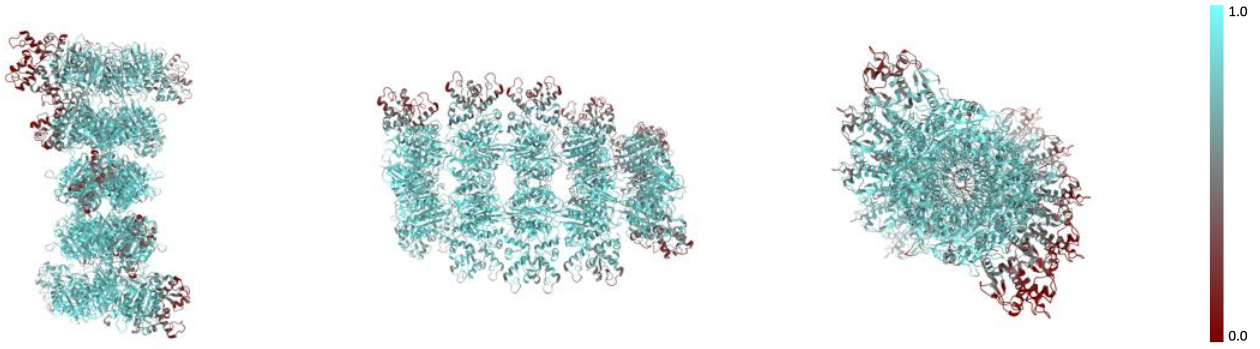
9.2 Q-score mapped to coordinate model (i)



X Y Z

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.

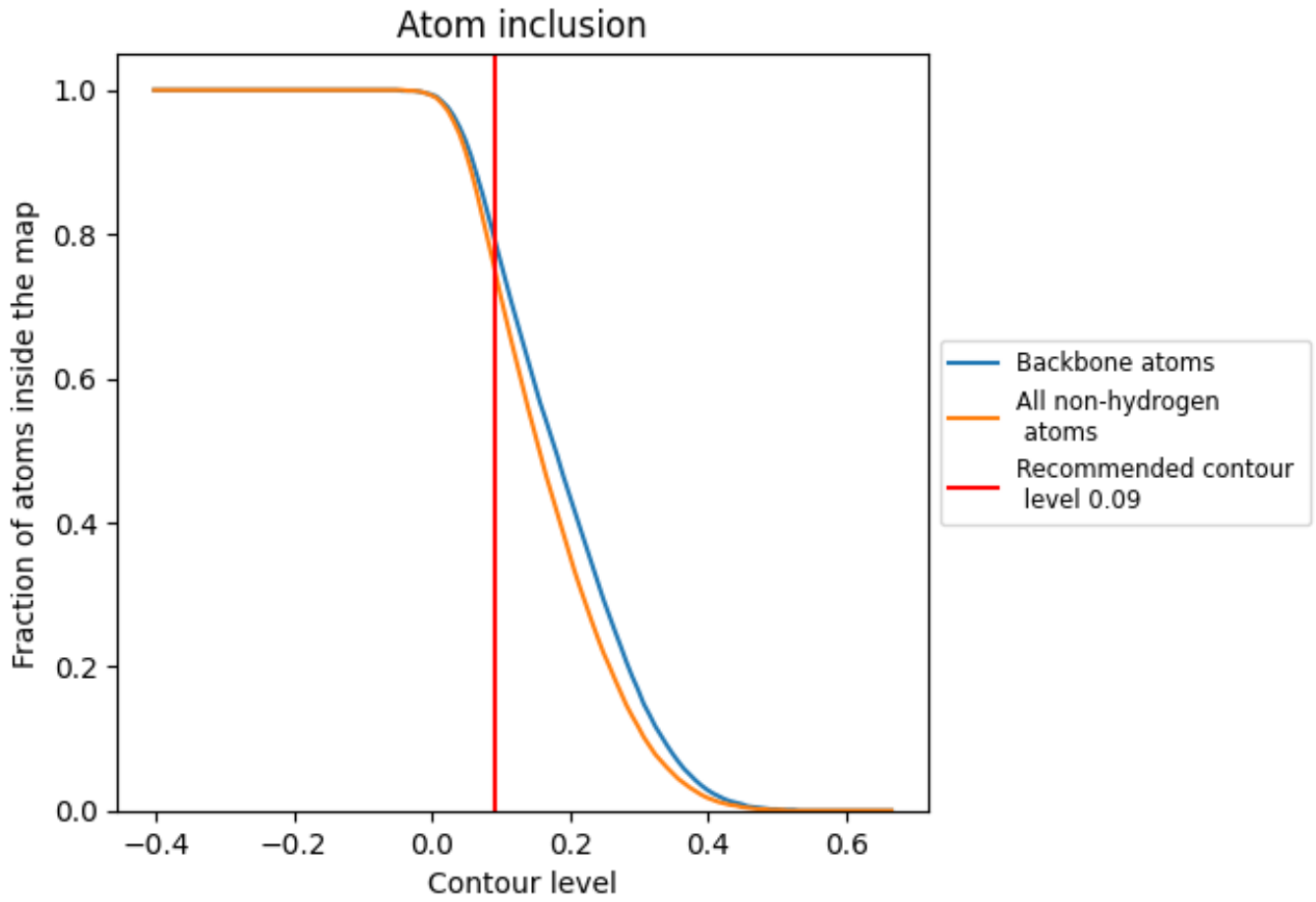
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.09).

9.4 Atom inclusion (i)



At the recommended contour level, 80% of all backbone atoms, 75% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.09) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7540	0.5450
A	0.8620	0.5670
B	0.8740	0.5720
C	0.7650	0.5460
D	0.7650	0.5430
E	0.8480	0.5610
F	0.7990	0.5550
G	0.8260	0.5610
H	0.6820	0.5320
I	0.8600	0.5660
J	0.8110	0.5600
K	0.8260	0.5630
L	0.7050	0.5440
M	0.7470	0.5400
N	0.6890	0.5320
O	0.7290	0.5370
P	0.5530	0.5060
Q	0.7560	0.5390
R	0.6930	0.5300
S	0.7120	0.5350
T	0.5760	0.5090

