



wwPDB EM Validation Summary Report ⓘ

Aug 31, 2024 – 01:24 PM EDT

PDB ID : 8UI0
EMDB ID : EMD-42285
Title : Structure of poised transcription complex Pol II-DSIF-NELF - pre-translocated
Authors : Vos, S.M.; Su, B.G.
Deposited on : 2023-10-09
Resolution : 2.70 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:

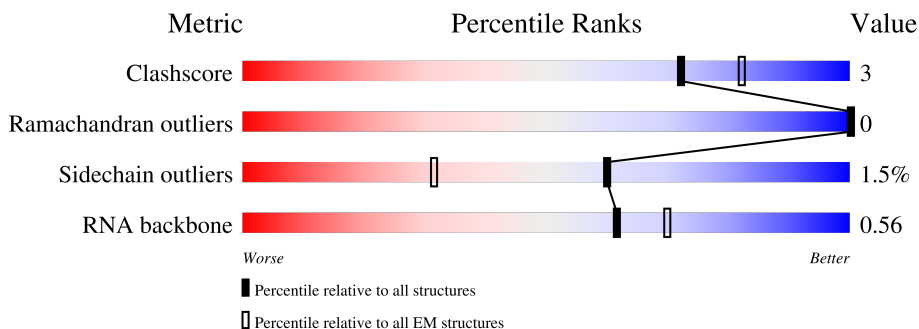
EMDB validation analysis : 0.0.1.dev112
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.38.3

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 2.70 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.






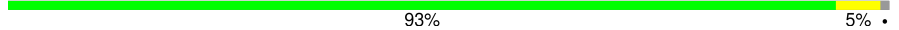


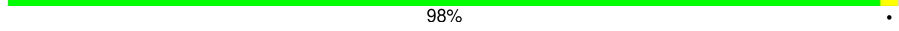


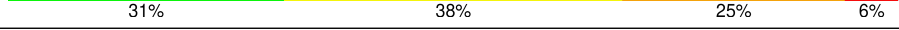



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415
RNA backbone	6643	2191

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	X	380	 96%
2	V	554	 55% 86% 13%
3	W	590	 26% 81% 8% 11%
4	A	1970	 64% 7% 29%
5	B	1174	 88% 9%
6	C	271	 88% 6% 6%
7	D	142	 76% 11% 13%

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Mol	Chain	Length	Quality of chain
8	E	210	 90% 9%
9	F	127	 57% 39%
10	G	171	 88% 11%
11	H	150	 93% 5%
12	I	125	 81% 10% 9%
13	J	67	 90% 9%
14	K	115	 98%
15	L	58	 64% 12% 24%
16	N	50	 52% 46%
17	P	16	 31% 38% 25% 6%
18	T	38	 89% 11%
19	U	527	 5% 30% 7% 63%
20	Z	1087	 6% 23% 75%

2 Entry composition i

There are 22 unique types of molecules in this entry. The entry contains 89334 atoms, of which 44428 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 Negative elongation factor E.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	H	N	O		
1	X	15	283	85	159	23	16	0	0

- Molecule 2 is a protein called Negative elongation factor B.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
2	V	550	8912	2825	4501	751	811	24	0	0

- Molecule 3 is a protein called Negative elongation factor C/D.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
3	W	526	8362	2673	4183	698	784	24	0	0

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
4	A	1399	22302	6976	11207	1985	2064	70	0	0

- Molecule 5 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
5	B	1136	18192	5739	9116	1597	1676	64	0	0

- Molecule 6 is a protein called DNA-directed RNA polymerase II subunit RPB3.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
6	C	254	4027	1282	1989	348	402	6	0	0

- Molecule 7 is a protein called RNA polymerase Rpb4/RPC9 core domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
7	D	124	2016	632	1004	173	203	4	0	0

- Molecule 8 is a protein called DNA-directed RNA polymerase II subunit E.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
8	E	207	3438	1083	1729	298	320	8	0	0

- Molecule 9 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC2.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
9	F	77	1271	397	653	105	111	5	0	0

- Molecule 10 is a protein called DNA-directed RNA polymerase subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
10	G	171	2710	875	1358	219	250	8	0	0

- Molecule 11 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC3.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
11	H	148	2333	750	1147	194	237	5	0	0

- Molecule 12 is a protein called DNA-directed RNA polymerase II subunit RPB9.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
12	I	114	1787	571	859	166	180	11	0	0

- Molecule 13 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC5.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
13	J	66	1064	339	540	88	91	6	0	0

- Molecule 14 is a protein called DNA-directed RNA polymerase II subunit RPB11-a.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
14	K	115	1863	593	942	152	174	2	0	0

- Molecule 15 is a protein called RNA polymerase II subunit K.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
15	L	44	751	231	378	72	64	6	0	0

- Molecule 16 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
16	N	27	856	263	305	97	164	27	0	0

- Molecule 17 is a RNA chain called RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
17	P	16	508	151	170	55	116	16	0	0

- Molecule 18 is a DNA chain called DNA (38-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
18	T	38	1209	371	427	148	225	38	0	0

- Molecule 19 is a protein called Negative elongation factor A.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
19	U	196	3054	953	1556	253	284	8	0	0

- Molecule 20 is a protein called Transcription elongation factor SPT5.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
20	Z	274	4387	1373	2205	393	405	11	0	0

- Molecule 21 is ZINC ION (three-letter code: ZN) (formula: Zn).

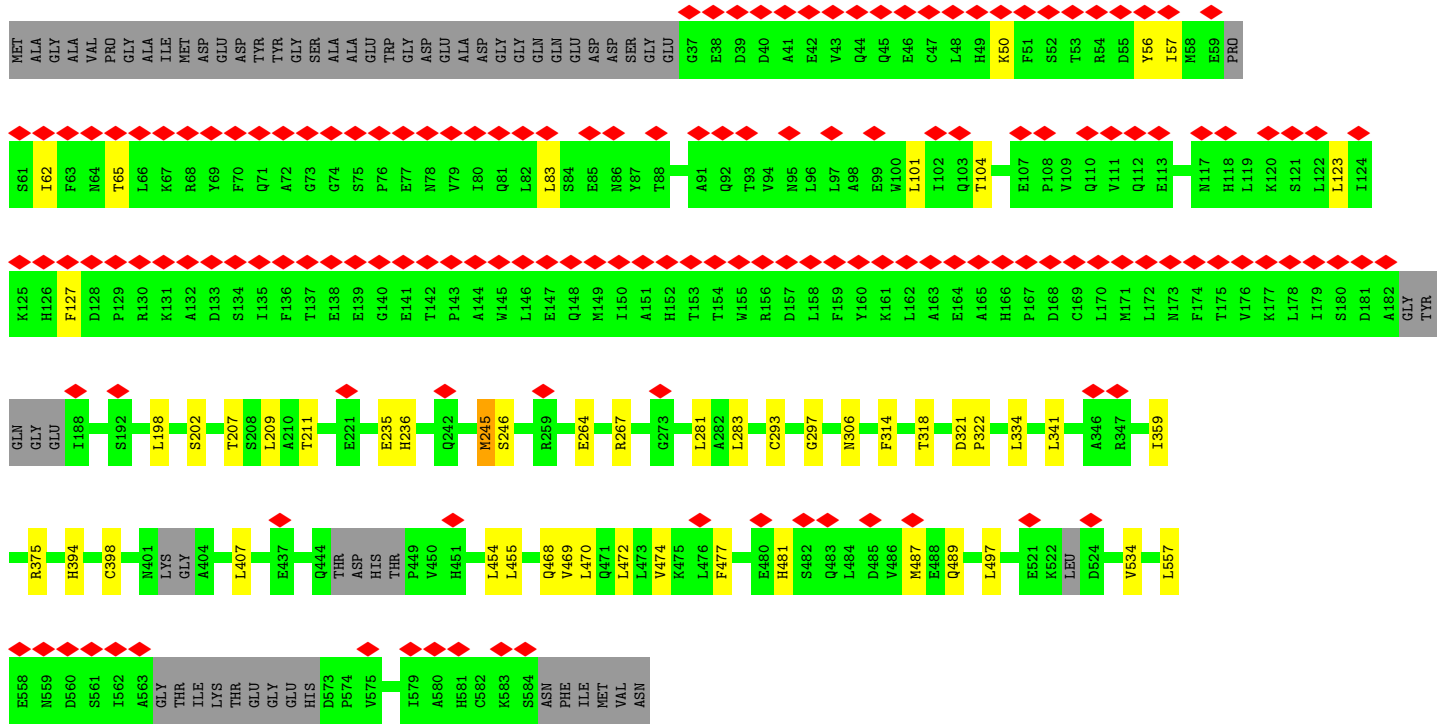
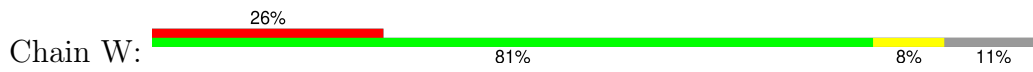
Mol	Chain	Residues	Atoms		AltConf
21	A	2	Total 2	Zn 2	0
21	B	1	Total 1	Zn 1	0
21	C	1	Total 1	Zn 1	0
21	I	2	Total 2	Zn 2	0
21	J	1	Total 1	Zn 1	0
21	L	1	Total 1	Zn 1	0

- Molecule 22 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
22	A	1	Total 1	Mg 1	0

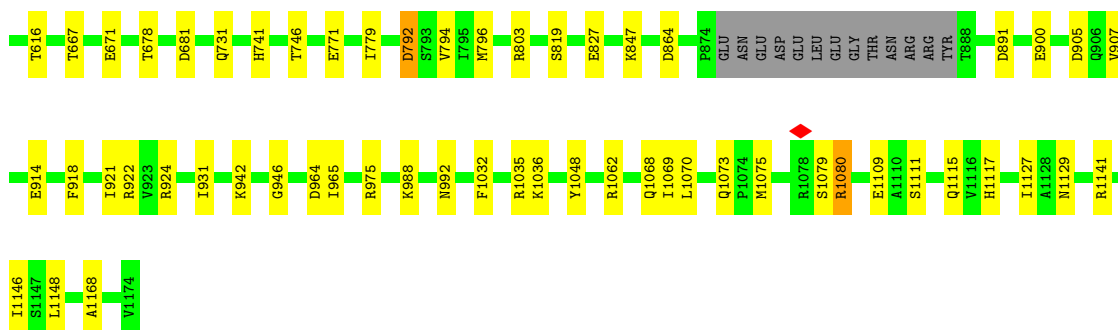


• Molecule 3: Negative elongation factor C/D



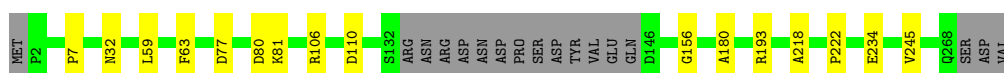
• Molecule 4: DNA-directed RNA polymerase subunit





- Molecule 6: DNA-directed RNA polymerase II subunit RPB3

Chain C: 88% 6% 6%



- Molecule 7: RNA polymerase Rpb4/RPC9 core domain-containing protein

Chain D: 76% 11% 13%



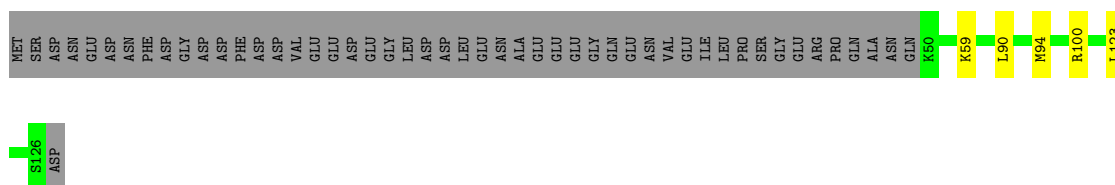
- Molecule 8: DNA-directed RNA polymerase II subunit E

Chain E: 90% 9%



- Molecule 9: DNA-directed RNA polymerases I, II, and III subunit RPABC2

Chain F: 57% 39%



- Molecule 10: DNA-directed RNA polymerase subunit

Chain G: 88% 11%




- Molecule 11: DNA-directed RNA polymerases I, II, and III subunit RPABC3

Chain H:  93% 5%




- Molecule 12: DNA-directed RNA polymerase II subunit RPB9

Chain I:  81% 10% 9%



- Molecule 13: DNA-directed RNA polymerases I, II, and III subunit RPABC5

Chain J:  90% 9%



- Molecule 14: DNA-directed RNA polymerase II subunit RPB11-a

Chain K:  98%



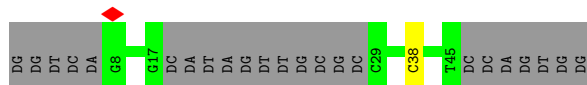
- Molecule 15: RNA polymerase II subunit K

Chain L:  64% 12% 24%



- Molecule 16: DNA

Chain N:  52% 46%

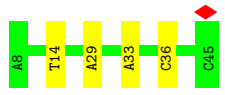
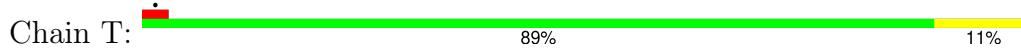


- Molecule 17: RNA

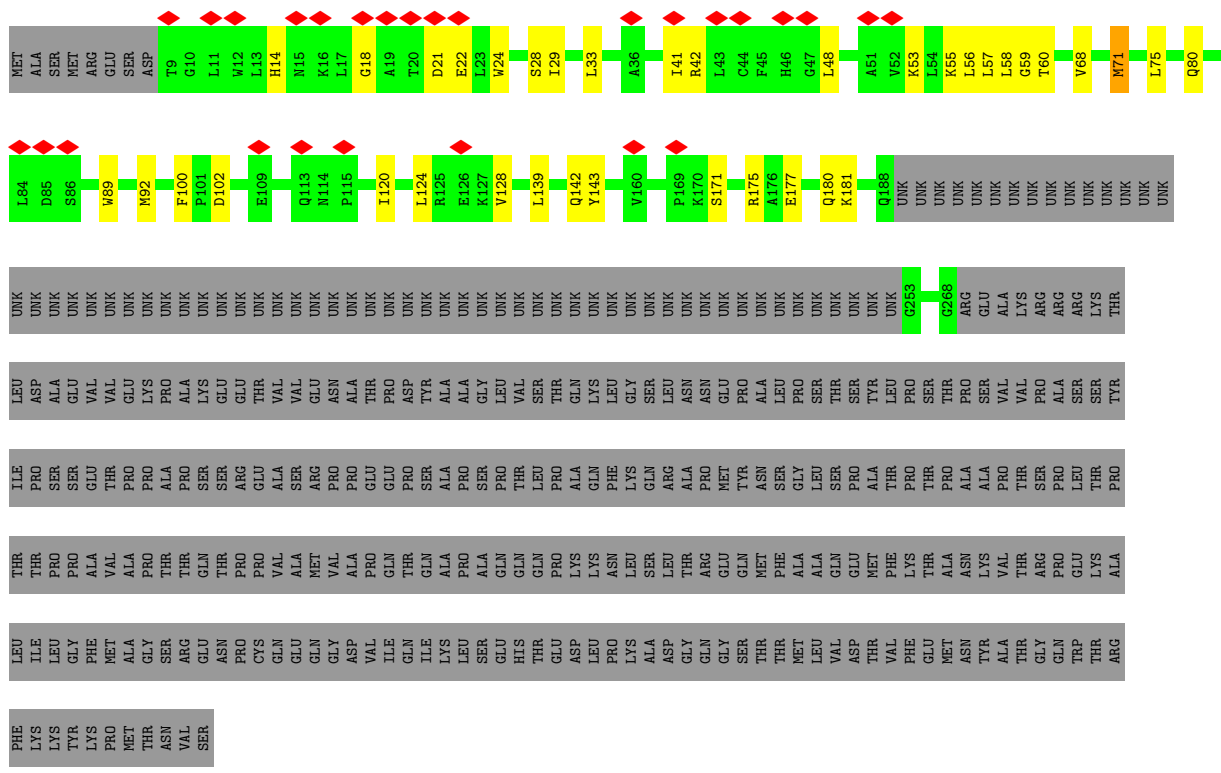
Chain P:  31% 38% 25% 6%



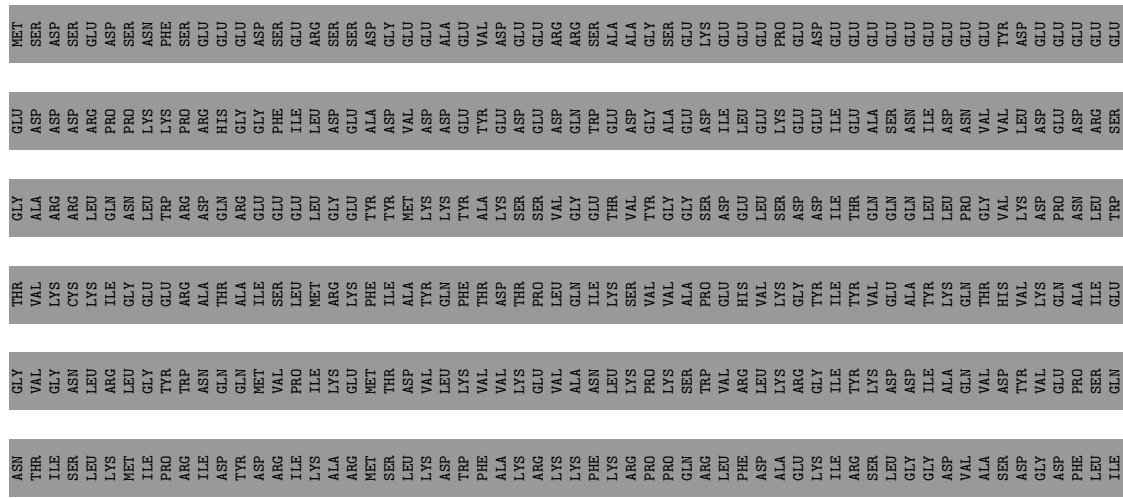
- Molecule 18: DNA (38-MER)



• Molecule 19: Negative elongation factor A



• Molecule 20: Transcription elongation factor SPT5



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	90283	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	51	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.739	Depositor
Minimum map value	-0.921	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.085	Depositor
Recommended contour level	0.35	Depositor
Map size (\AA)	420.864, 420.864, 420.864	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.822, 0.822, 0.822	Depositor

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MG

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	X	0.37	0/124	0.60	0/159
2	V	0.34	0/4496	0.69	3/6074 (0.0%)
3	W	0.31	0/4261	0.60	0/5777
4	A	0.30	0/11294	0.67	2/15241 (0.0%)
5	B	0.30	0/9257	0.65	0/12493
6	C	0.29	0/2081	0.64	1/2828 (0.0%)
7	D	0.30	0/1025	0.61	0/1376
8	E	0.32	0/1739	0.70	1/2347 (0.0%)
9	F	0.30	0/628	0.67	0/848
10	G	0.33	0/1383	0.64	0/1874
11	H	0.30	0/1207	0.66	0/1628
12	I	0.34	0/949	0.73	1/1284 (0.1%)
13	J	0.34	0/533	0.69	0/719
14	K	0.31	0/940	0.59	0/1271
15	L	0.37	0/378	0.86	1/500 (0.2%)
16	N	0.85	0/615	1.08	0/944
17	P	0.34	0/376	1.15	4/583 (0.7%)
18	T	0.81	1/878 (0.1%)	1.06	2/1353 (0.1%)
19	U	0.31	0/1521	0.63	0/2063
20	Z	0.27	0/2214	0.63	1/2979 (0.0%)
All	All	0.34	1/45899 (0.0%)	0.68	16/62341 (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
4	A	0	1
5	B	0	1
12	I	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
19	U	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	T	36	DC	C1'-N1	6.83	1.58	1.49

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	110	ASP	CB-CG-OD1	9.24	126.61	118.30
4	A	849	ASP	CB-CG-OD1	8.00	125.50	118.30
15	L	56	ASP	CB-CG-OD2	7.89	125.41	118.30
2	V	221	PHE	CB-CG-CD2	7.67	126.17	120.80
4	A	1423	ASP	CB-CG-OD2	7.29	124.86	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	A	532	ARG	Sidechain
5	B	242	ARG	Sidechain
12	I	103	ARG	Sidechain
19	U	42	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	X	124	159	159	0	0
2	V	4411	4501	4501	46	0
3	W	4179	4183	4181	33	0
4	A	11095	11207	11207	81	0
5	B	9076	9116	9116	62	0
6	C	2038	1989	1989	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	D	1012	1004	1004	10	0
8	E	1709	1729	1729	12	0
9	F	618	653	653	3	0
10	G	1352	1358	1358	11	0
11	H	1186	1147	1147	5	0
12	I	928	859	859	8	0
13	J	524	540	540	5	0
14	K	921	942	942	2	0
15	L	373	378	378	4	0
16	N	551	305	307	1	0
17	P	338	170	170	5	0
18	T	782	427	427	2	0
19	U	1498	1556	1556	21	0
20	Z	2182	2205	2205	10	0
21	A	2	0	0	0	0
21	B	1	0	0	0	0
21	C	1	0	0	0	0
21	I	2	0	0	0	0
21	J	1	0	0	0	0
21	L	1	0	0	0	0
22	A	1	0	0	0	0
All	All	44906	44428	44428	287	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 287 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)
5:B:237:VAL:HG21	5:B:369:VAL:HG12	1.57	0.86
19:U:57:LEU:O	19:U:60:THR:OG1	2.06	0.74
7:D:73:ARG:NE	7:D:102:ASN:O	2.23	0.72
2:V:79:ALA:O	2:V:83:GLU:N	2.23	0.70
5:B:864:ASP:OD1	20:Z:725:LYS:NZ	2.25	0.69

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	X	13/380 (3%)	13 (100%)	0	0	100	100
2	V	548/554 (99%)	530 (97%)	18 (3%)	0	100	100
3	W	512/590 (87%)	493 (96%)	19 (4%)	0	100	100
4	A	1387/1970 (70%)	1337 (96%)	50 (4%)	0	100	100
5	B	1130/1174 (96%)	1076 (95%)	54 (5%)	0	100	100
6	C	250/271 (92%)	242 (97%)	8 (3%)	0	100	100
7	D	122/142 (86%)	115 (94%)	7 (6%)	0	100	100
8	E	203/210 (97%)	196 (97%)	7 (3%)	0	100	100
9	F	75/127 (59%)	73 (97%)	2 (3%)	0	100	100
10	G	169/171 (99%)	165 (98%)	4 (2%)	0	100	100
11	H	146/150 (97%)	143 (98%)	3 (2%)	0	100	100
12	I	112/125 (90%)	104 (93%)	8 (7%)	0	100	100
13	J	64/67 (96%)	62 (97%)	2 (3%)	0	100	100
14	K	113/115 (98%)	111 (98%)	2 (2%)	0	100	100
15	L	42/58 (72%)	41 (98%)	1 (2%)	0	100	100
19	U	192/527 (36%)	189 (98%)	3 (2%)	0	100	100
20	Z	266/1087 (24%)	256 (96%)	10 (4%)	0	100	100
All	All	5344/7718 (69%)	5146 (96%)	198 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	X	13/331 (4%)	13 (100%)	0	100	100
2	V	493/495 (100%)	486 (99%)	7 (1%)	62	84
3	W	466/513 (91%)	463 (99%)	3 (1%)	84	94
4	A	1232/1749 (70%)	1215 (99%)	17 (1%)	62	84
5	B	994/1027 (97%)	975 (98%)	19 (2%)	52	79
6	C	231/248 (93%)	231 (100%)	0	100	100
7	D	114/126 (90%)	111 (97%)	3 (3%)	41	70
8	E	190/192 (99%)	187 (98%)	3 (2%)	58	82
9	F	67/111 (60%)	65 (97%)	2 (3%)	36	65
10	G	152/152 (100%)	147 (97%)	5 (3%)	33	62
11	H	129/131 (98%)	128 (99%)	1 (1%)	79	91
12	I	103/112 (92%)	103 (100%)	0	100	100
13	J	55/56 (98%)	54 (98%)	1 (2%)	54	80
14	K	104/104 (100%)	104 (100%)	0	100	100
15	L	41/55 (74%)	40 (98%)	1 (2%)	44	73
19	U	167/396 (42%)	161 (96%)	6 (4%)	30	59
20	Z	244/940 (26%)	240 (98%)	4 (2%)	58	82
All	All	4795/6738 (71%)	4723 (98%)	72 (2%)	60	83

5 of 72 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
10	G	131	MET
20	Z	740	CYS
11	H	145	MET
19	U	100	PHE
4	A	1428	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
17	P	16/16 (100%)	9 (56%)	2 (12%)

5 of 9 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
17	P	4	U
17	P	5	U
17	P	6	U
17	P	7	G
17	P	8	G

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
17	P	3	U
17	P	6	U

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](#)

Of 9 ligands modelled in this entry, 9 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

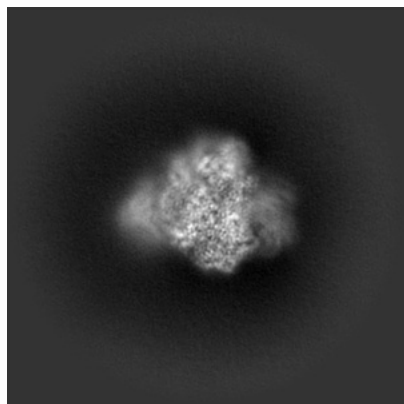
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-42285. These allow visual inspection of the internal detail of the map and identification of artifacts.

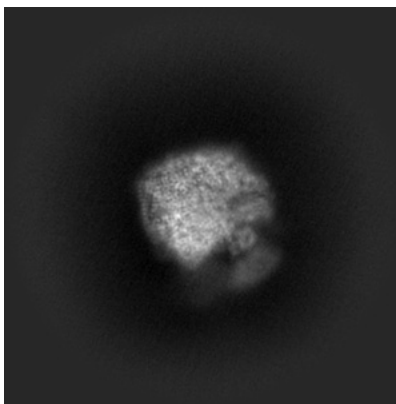
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

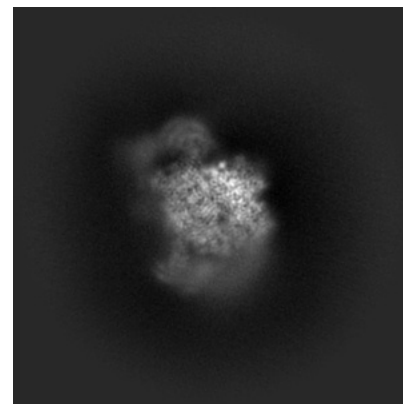
6.1.1 Primary map



X

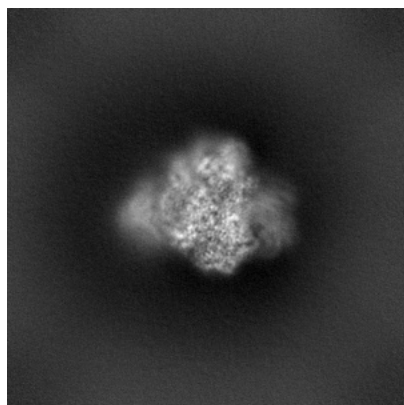


Y

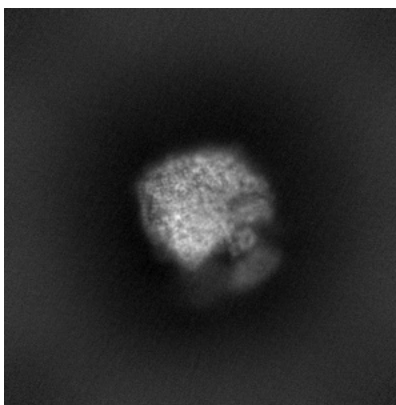


Z

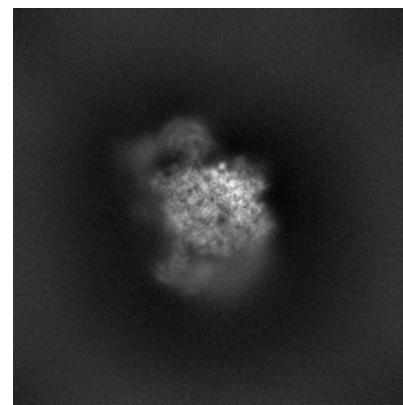
6.1.2 Raw map



X



Y

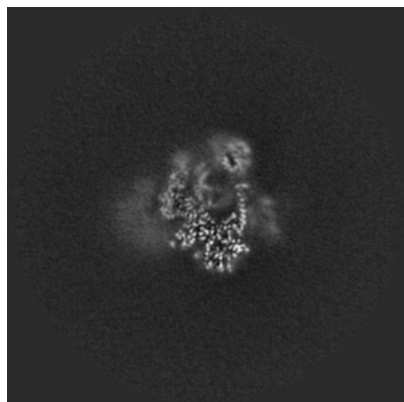


Z

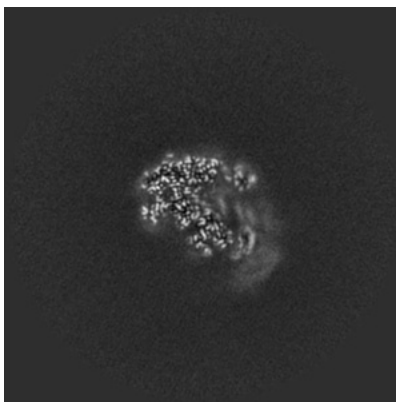
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

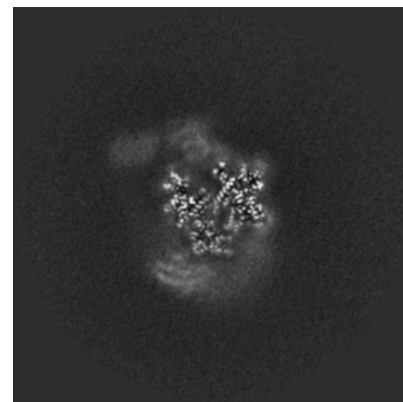
6.2.1 Primary map



X Index: 256

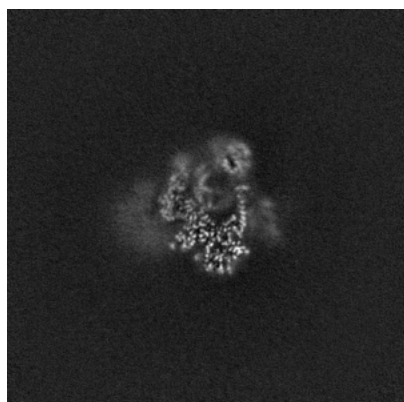


Y Index: 256

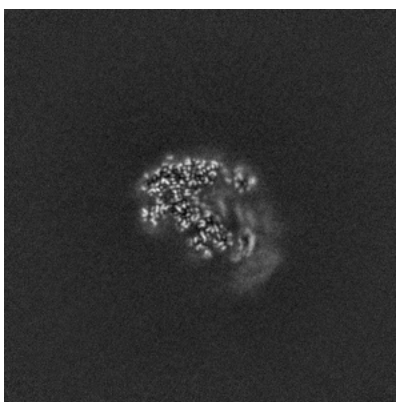


Z Index: 256

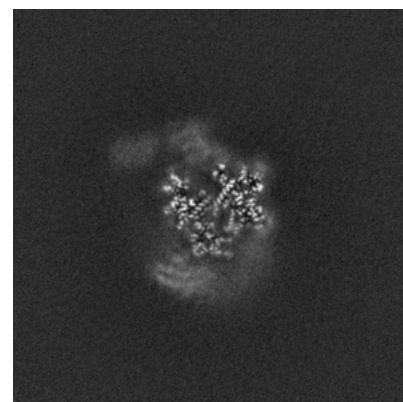
6.2.2 Raw map



X Index: 256



Y Index: 256

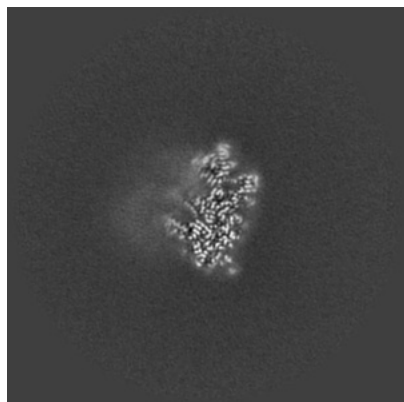


Z Index: 256

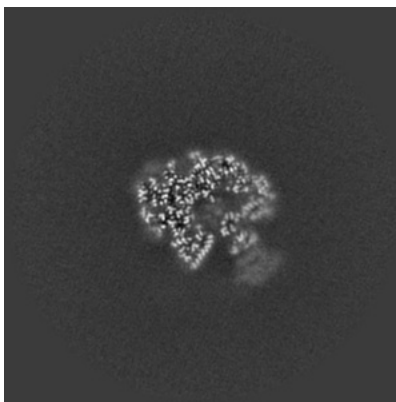
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

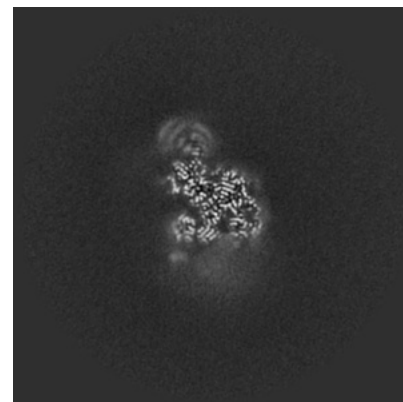
6.3.1 Primary map



X Index: 289

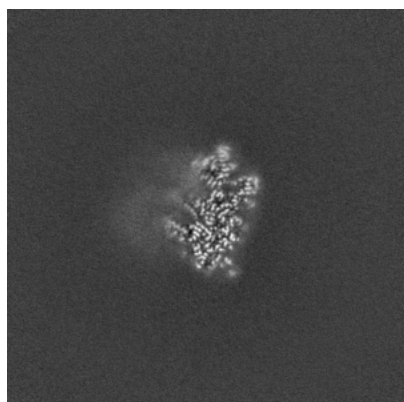


Y Index: 281

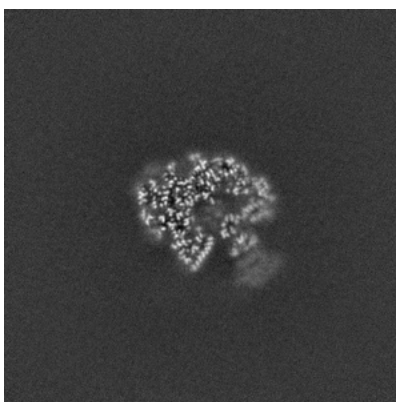


Z Index: 220

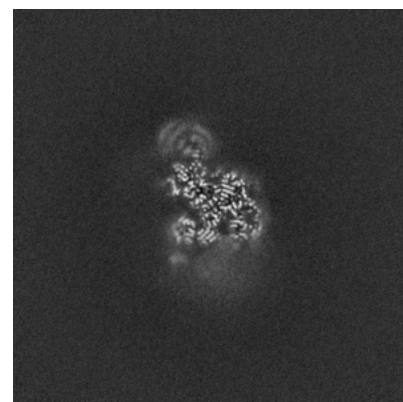
6.3.2 Raw map



X Index: 289



Y Index: 281

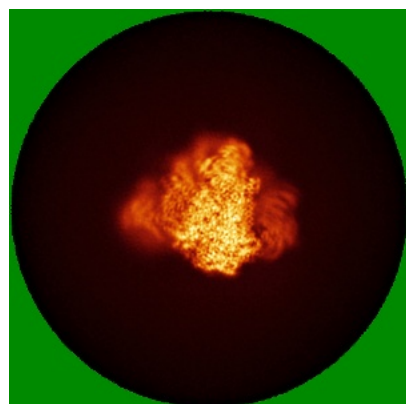


Z Index: 220

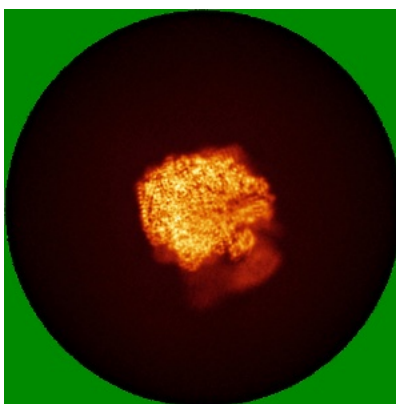
The images above show the largest variance slices of the map in three orthogonal directions.

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

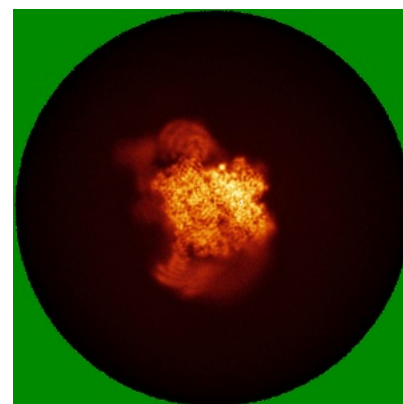
6.4.1 Primary map



X

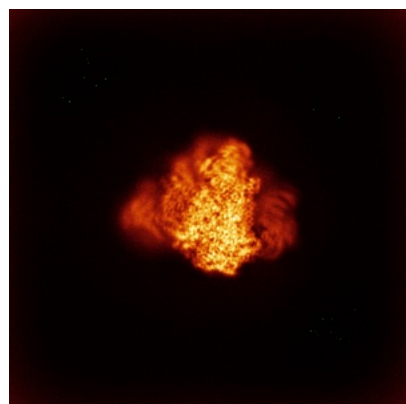


Y

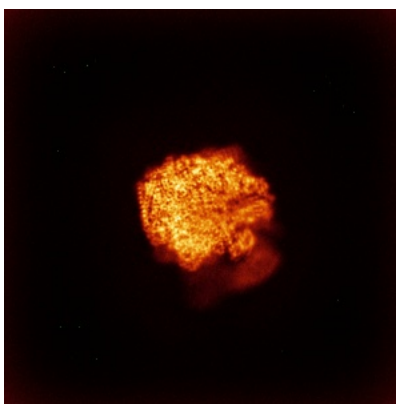


Z

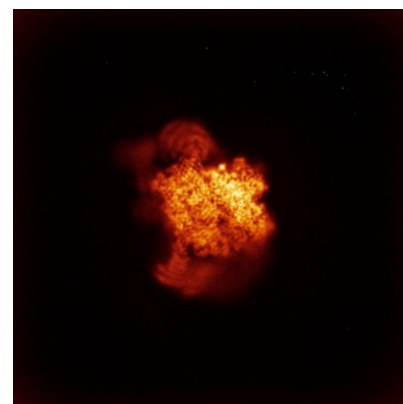
6.4.2 Raw map



X



Y

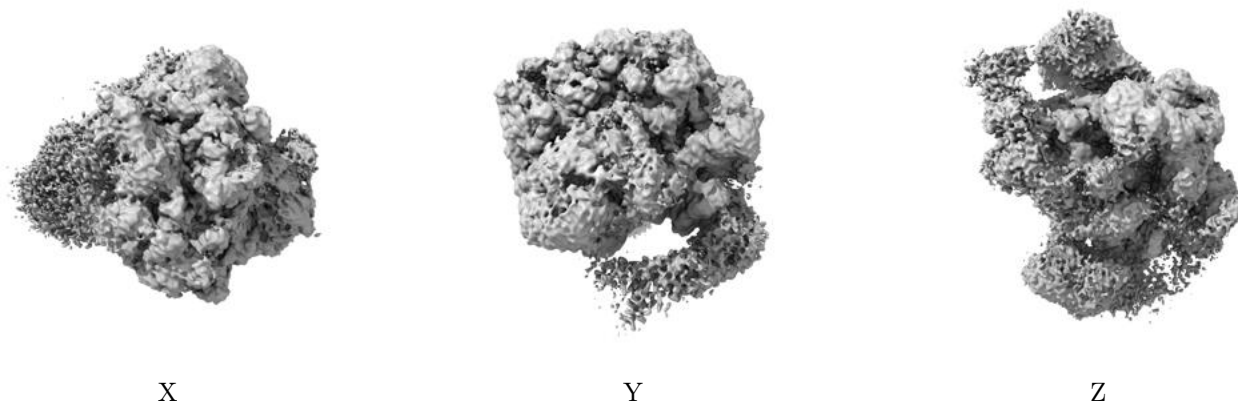


Z

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.

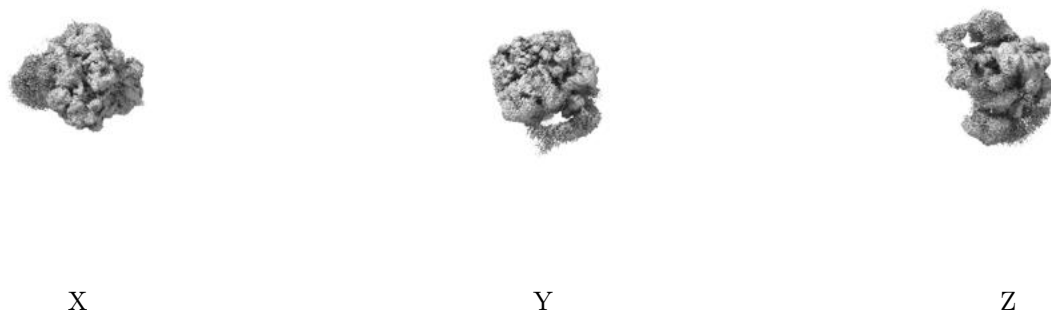
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.35. 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



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.

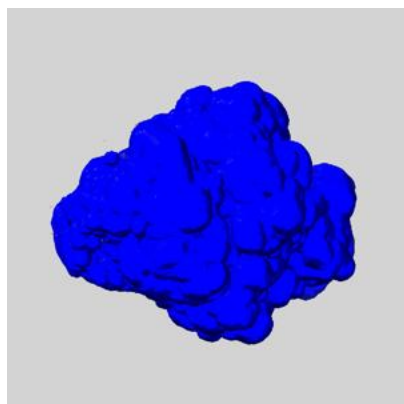
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

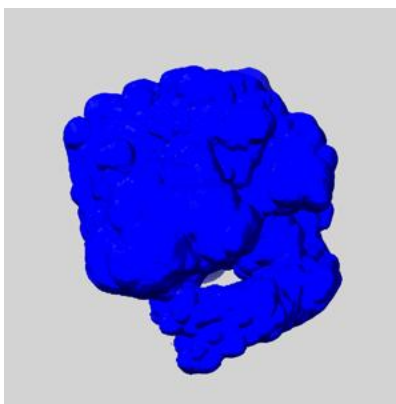
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

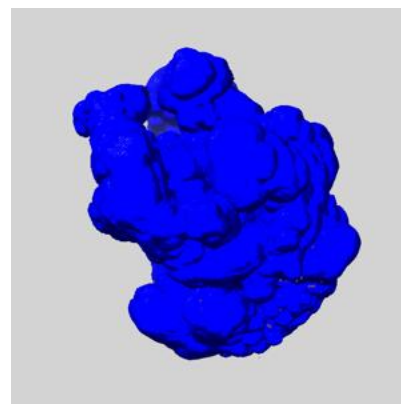
6.6.1 emd_42285_msk_1.map [i](#)



X



Y

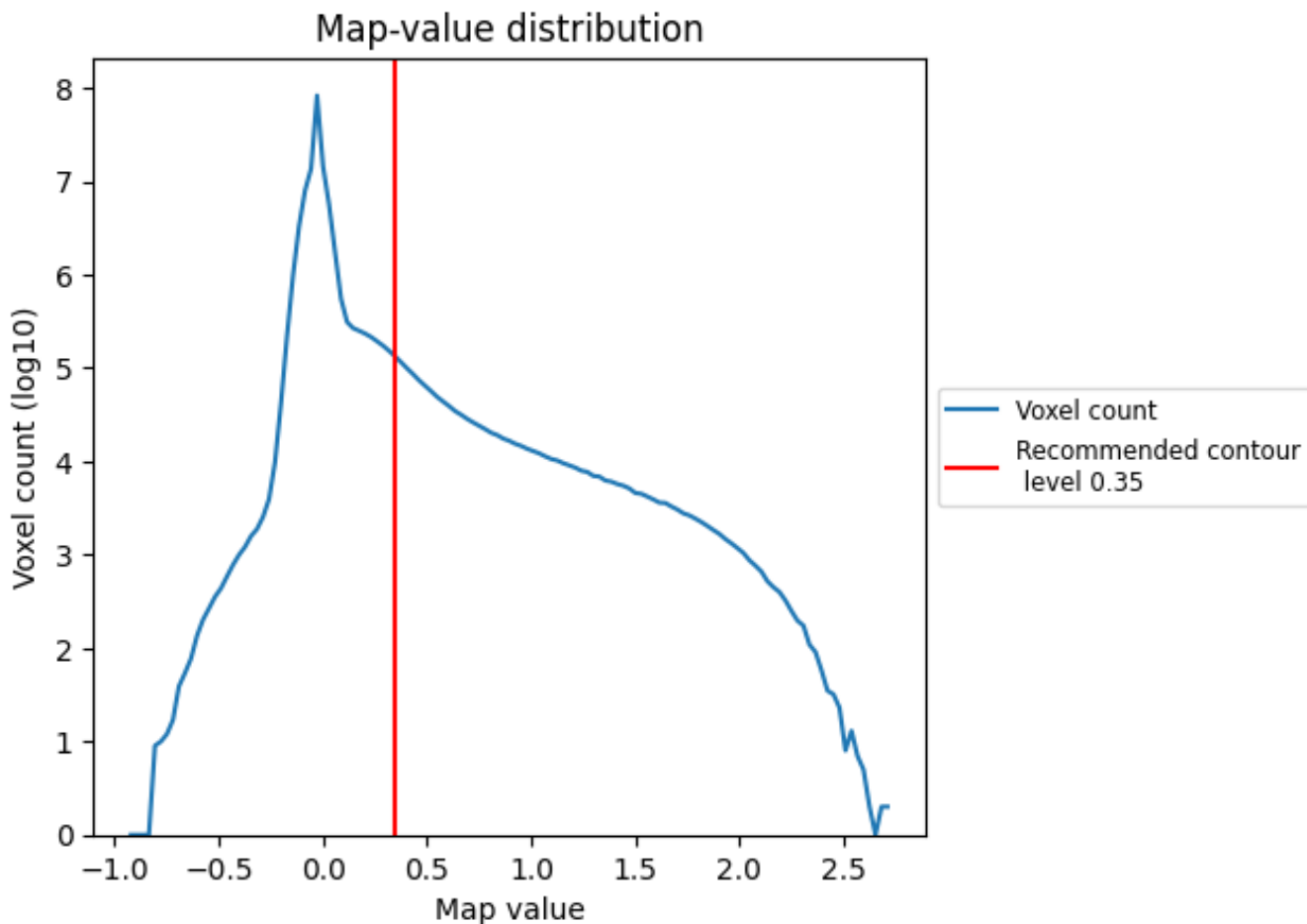


Z

7 Map analysis [i](#)

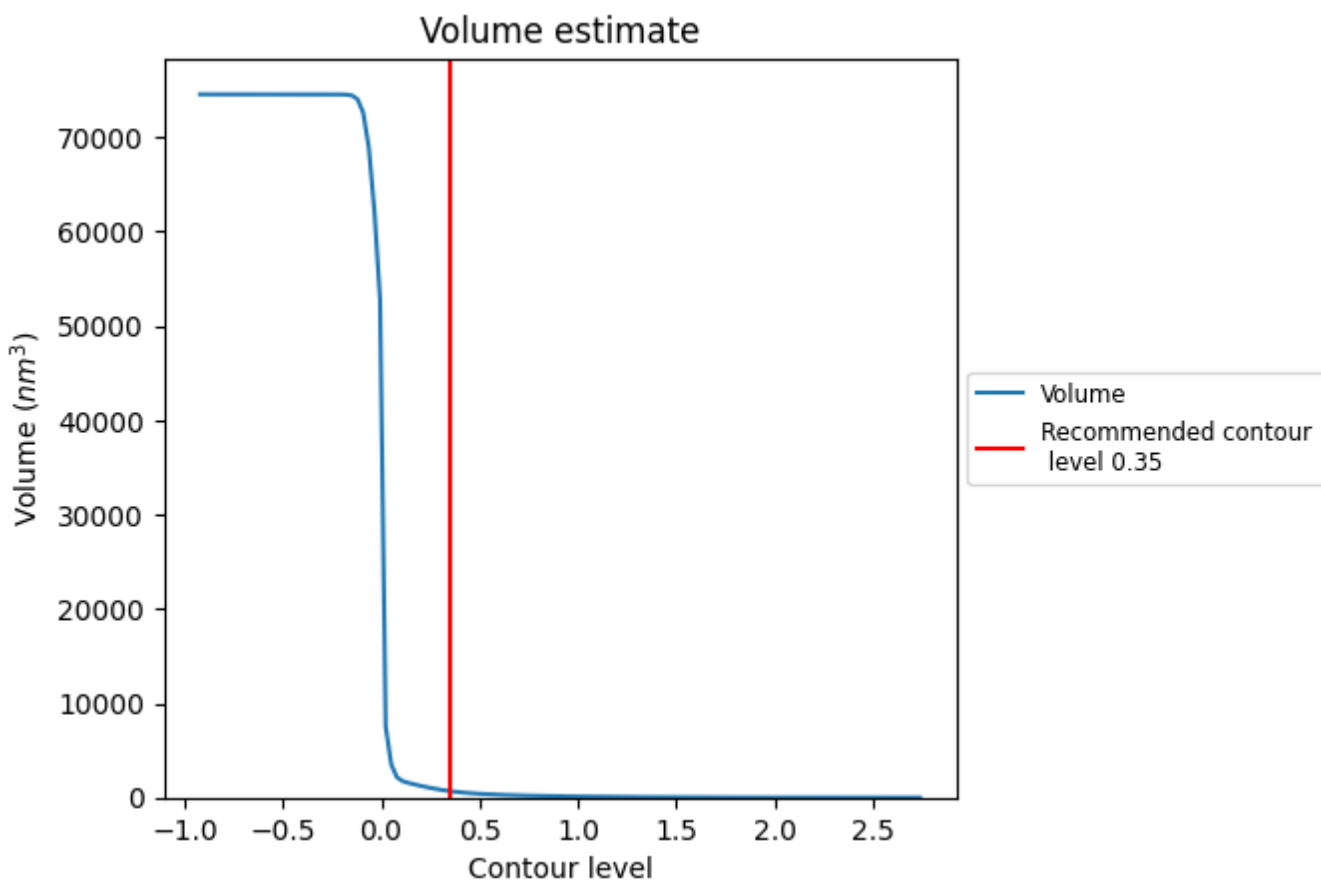
This section contains the results of statistical analysis of the map.

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.

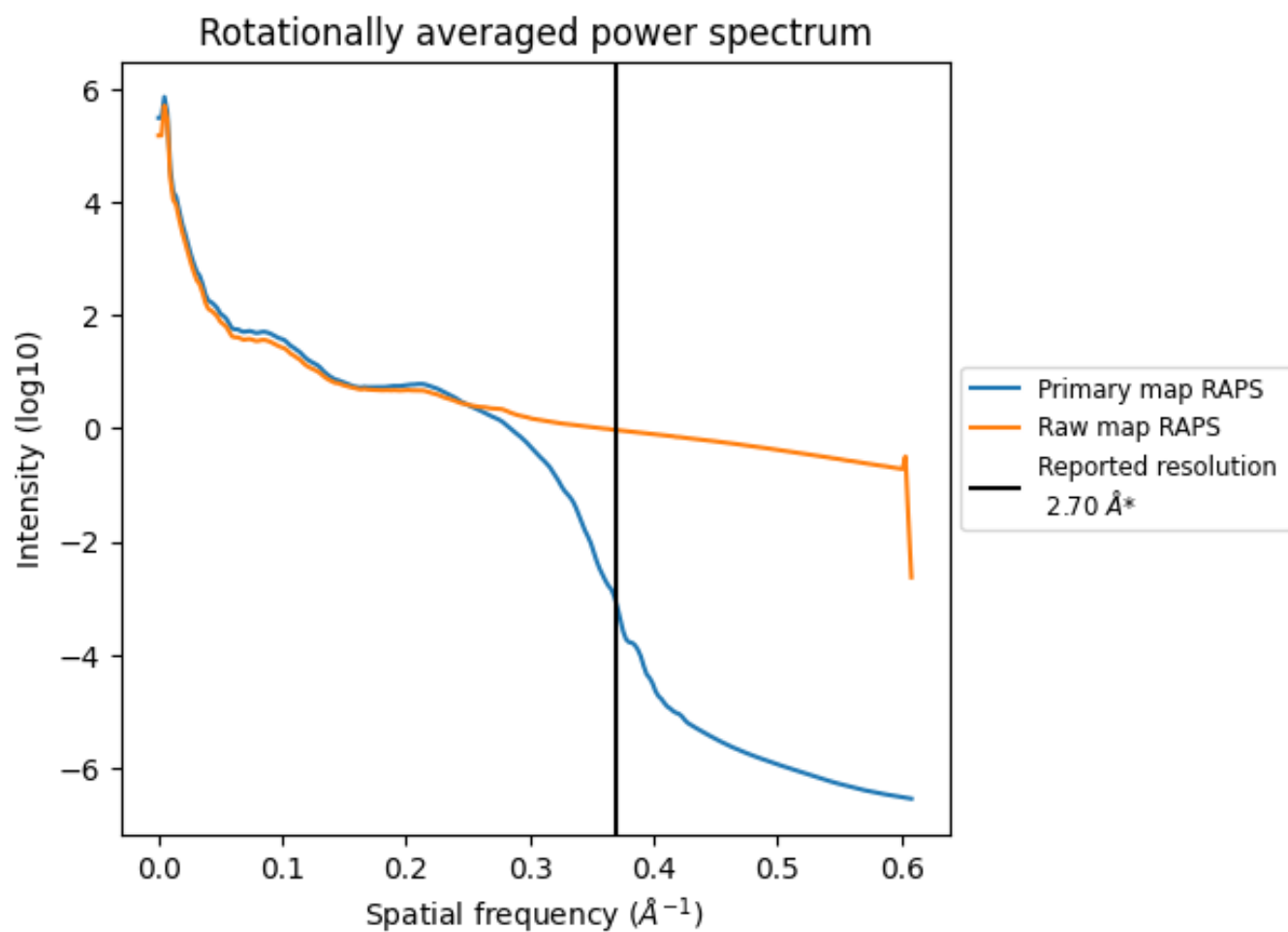
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 683 nm³; this corresponds to an approximate mass of 617 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

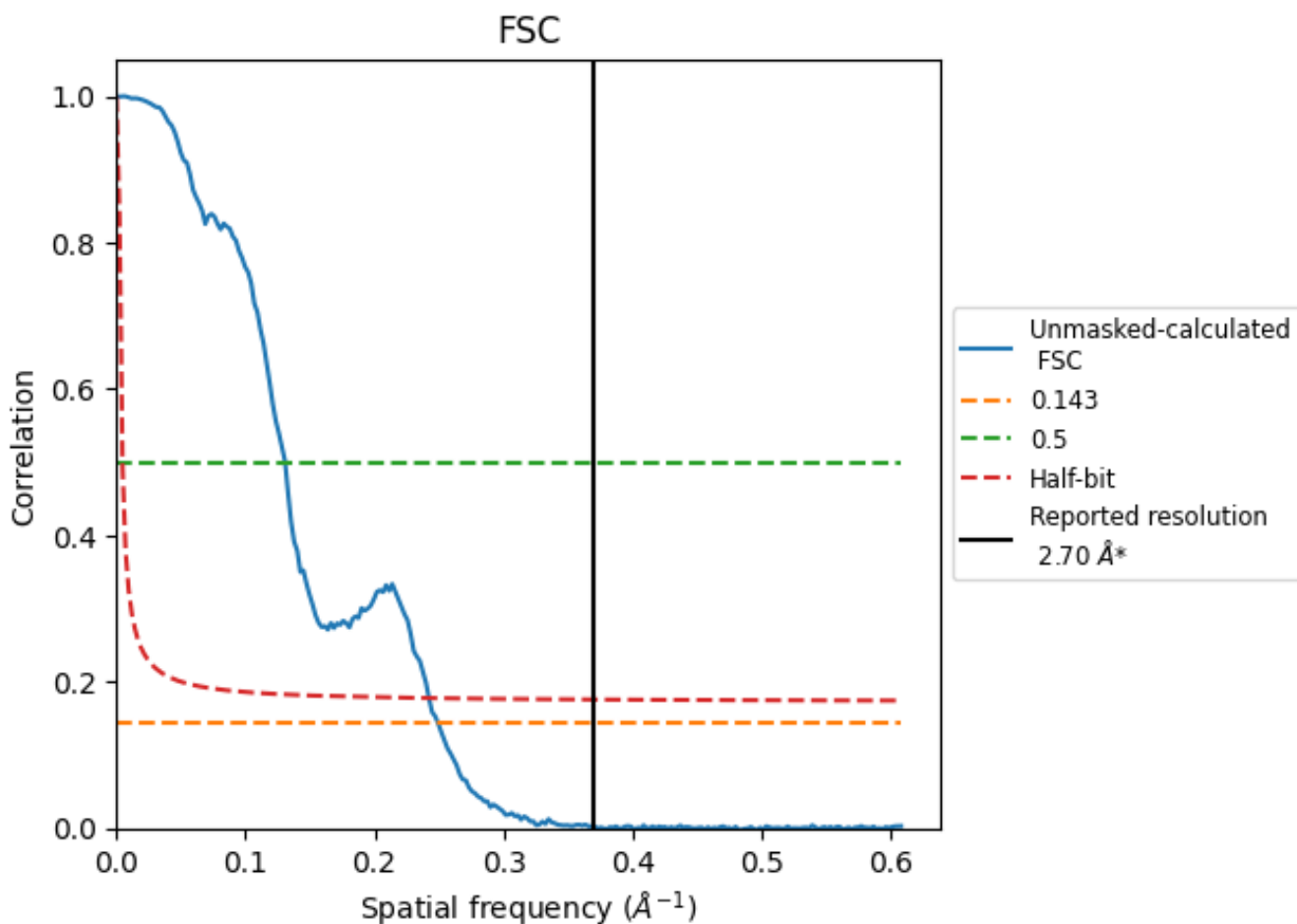


*Reported resolution corresponds to spatial frequency of 0.370 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.370 Å⁻¹

8.2 Resolution estimates [i](#)

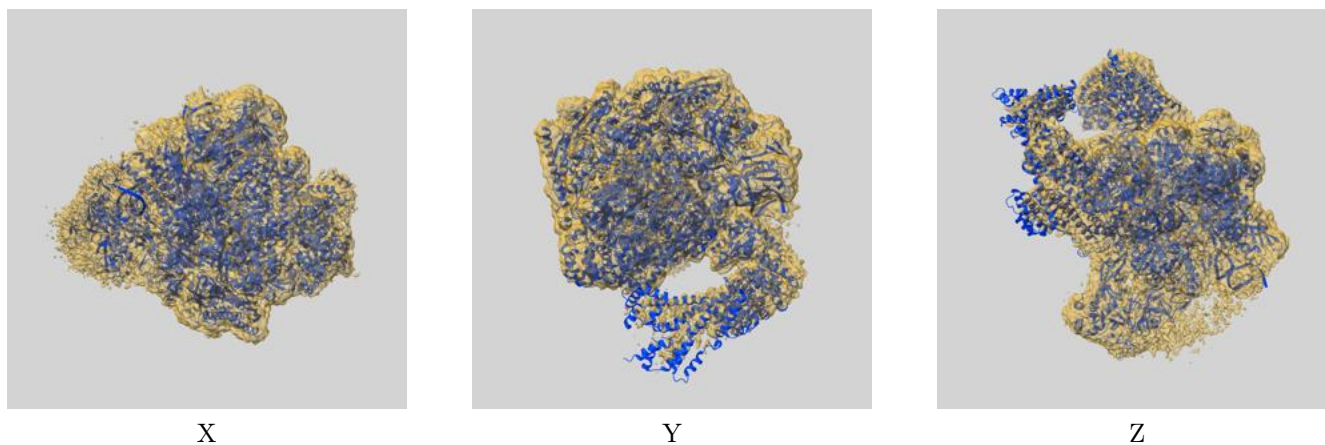
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.70	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.01	7.65	4.13

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.01 differs from the reported value 2.7 by more than 10 %

9 Map-model fit [i](#)

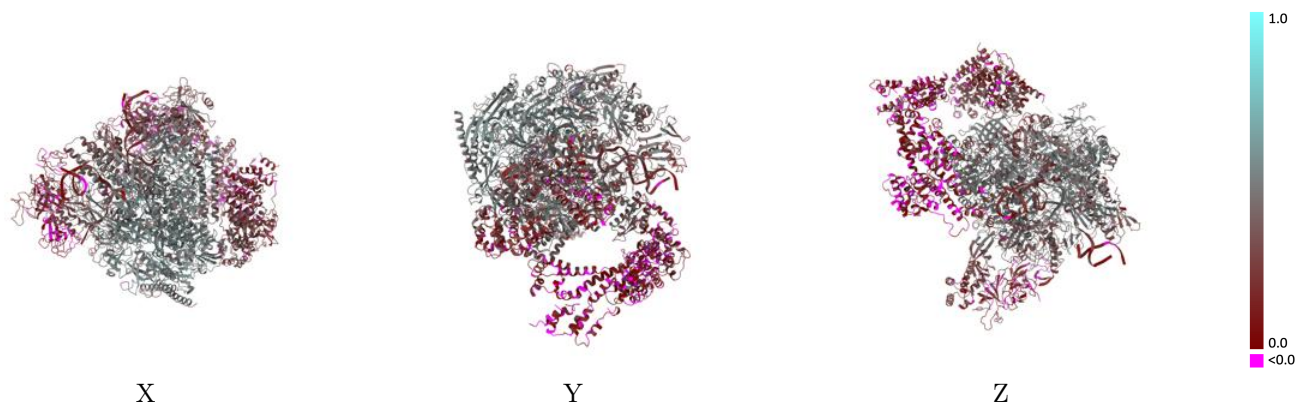
This section contains information regarding the fit between EMDB map EMD-42285 and PDB model 8UI0. Per-residue inclusion information can be found in section 3 on page 8.

9.1 Map-model overlay [i](#)



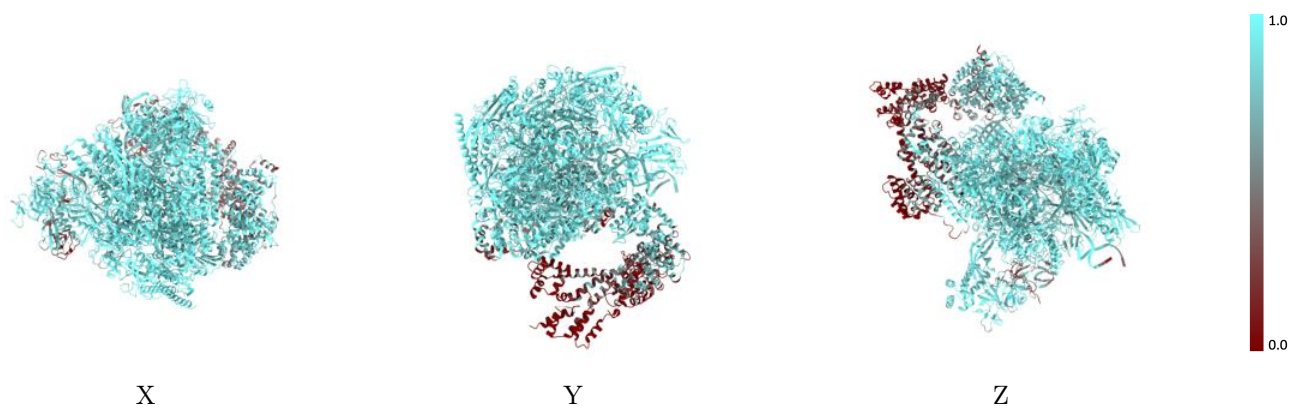
The images above show the 3D surface view of the map at the recommended contour level 0.35 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



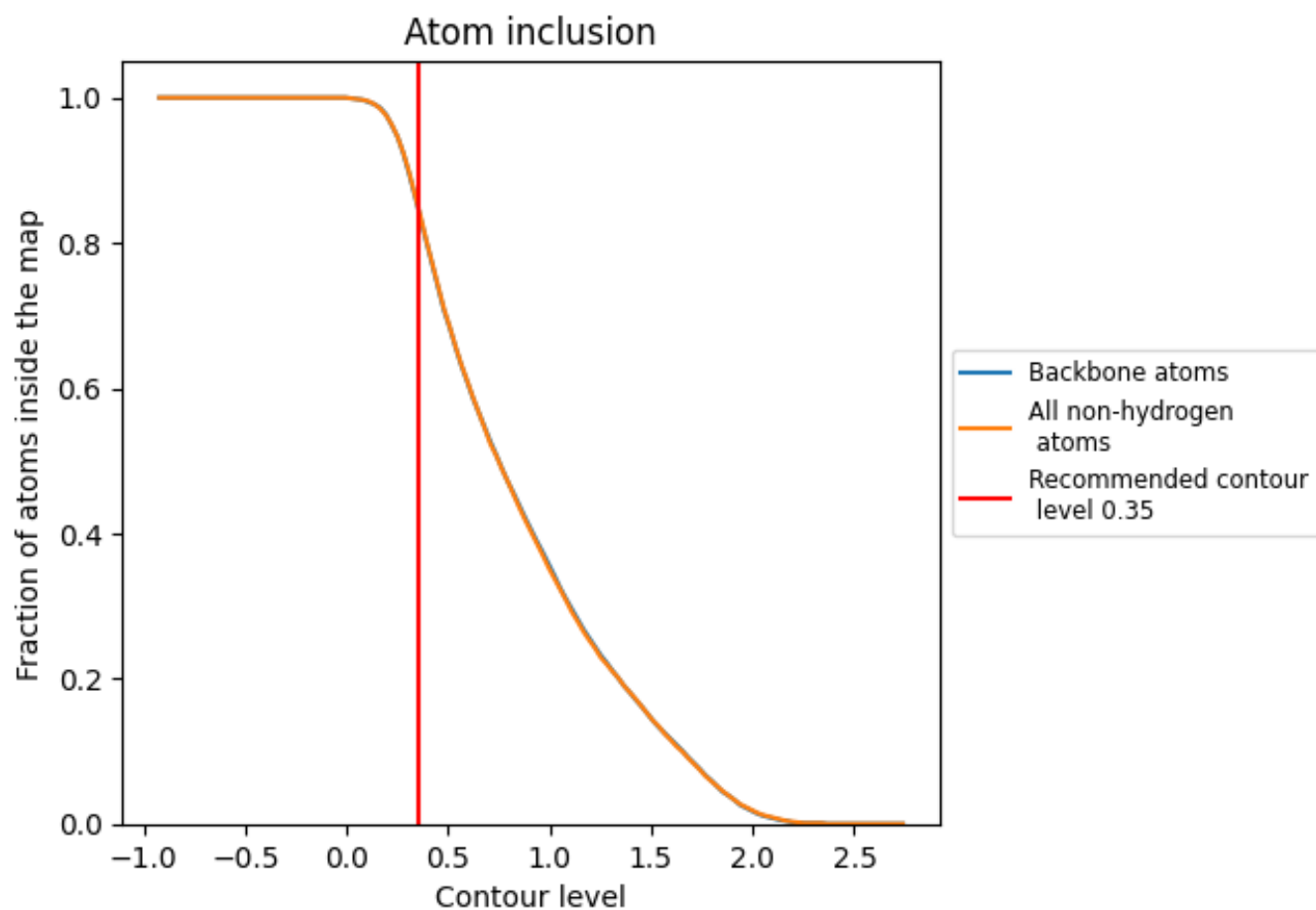
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.35).











































9.4 Atom inclusion [i](#)



At the recommended contour level, 85% of all backbone atoms, 85% 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 (0.35) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8530	 0.3530
A	 0.9590	 0.4460
B	 0.9690	 0.4730
C	 0.9850	 0.5010
D	 0.9730	 0.2110
E	 0.9640	 0.4120
F	 0.9900	 0.4900
G	 0.9690	 0.2600
H	 0.9760	 0.4830
I	 0.9880	 0.4140
J	 0.9840	 0.5190
K	 0.9800	 0.5000
L	 0.9920	 0.4680
N	 0.8710	 0.1500
P	 0.9320	 0.3250
T	 0.9120	 0.2540
U	 0.7450	 0.1880
V	 0.3900	 0.0810
W	 0.6290	 0.1790
X	 0.7900	 0.1060
Z	 0.6690	 0.1580

