



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 02:51 AM UTC

PDB ID : 8W7S / pdb_00008w7s
EMDB ID : EMD-37345
Title : Yeast replisome in state IV
Authors : Dang, S.; Zhai, Y.; Feng, J.; Yu, D.
Deposited on : 2023-08-31
Resolution : 7.39 Å(reported)
Based on initial model : 6SKL

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 ⓘ symbol.

The types of validation reports are described at

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

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

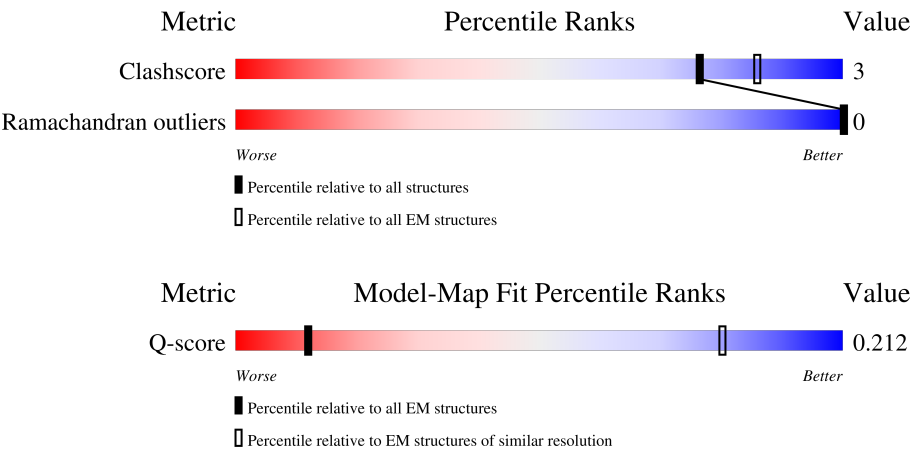
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 7.39 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	-
Q-score	-	25397	390 (6.89 - 7.86)

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	2	868	22% 69% 29%
2	3	971	16% 61% 37%
3	4	933	23% 65% 33%
4	5	775	9% 65% 33%
5	6	1017	22% 59% 40%

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Mol	Chain	Length	Quality of chain
6	7	845	
7	A	208	
8	B	213	
9	C	194	
10	D	294	
11	E	650	
12	F	927	
12	G	927	
12	H	927	
13	I	71	
14	N	689	

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 25521 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 DNA replication licensing factor MCM2.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	2	617	Total	C	N	O	0	0
			2469	1234	617	618		

- Molecule 2 is a protein called DNA replication licensing factor MCM3.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	3	612	Total	C	N	O	0	0
			2448	1224	612	612		

- Molecule 3 is a protein called DNA replication licensing factor MCM4.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	4	623	Total	C	N	O	0	0
			2492	1246	623	623		

- Molecule 4 is a protein called Minichromosome maintenance protein 5.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	5	521	Total	C	N	O	0	0
			2084	1042	521	521		

- Molecule 5 is a protein called DNA replication licensing factor MCM6.

Mol	Chain	Residues	Atoms				AltConf	Trace
5	6	613	Total	C	N	O	0	0
			2452	1226	613	613		

- Molecule 6 is a protein called DNA replication licensing factor MCM7.

Mol	Chain	Residues	Atoms				AltConf	Trace
6	7	633	Total	C	N	O	0	0
			2532	1266	633	633		

- Molecule 7 is a protein called DNA replication complex GINS protein PSF1.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	A	197	Total	C	N	O	0	0
			789	394	197	198		

- Molecule 8 is a protein called DNA replication complex GINS protein PSF2.

Mol	Chain	Residues	Atoms				AltConf	Trace
8	B	193	Total	C	N	O	0	0
			772	386	193	193		

- Molecule 9 is a protein called DNA replication complex GINS protein PSF3.

Mol	Chain	Residues	Atoms				AltConf	Trace
9	C	172	Total	C	N	O	0	0
			689	344	172	173		

- Molecule 10 is a protein called DNA replication complex GINS protein SLD5.

Mol	Chain	Residues	Atoms				AltConf	Trace
10	D	243	Total	C	N	O	0	0
			973	486	243	244		

- Molecule 11 is a protein called Cell division control protein 45.

Mol	Chain	Residues	Atoms				AltConf	Trace
11	E	564	Total	C	N	O	0	0
			2257	1128	564	565		

- Molecule 12 is a protein called DNA polymerase alpha-binding protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
12	F	424	Total	C	N	O	0	0
			1696	848	424	424		
12	G	422	Total	C	N	O	0	0
			1688	844	422	422		
12	H	425	Total	C	N	O	0	0
			1700	850	425	425		

- Molecule 13 is a DNA chain called DNA (71-mer).

Mol	Chain	Residues	Atoms					AltConf	Trace
13	I	7	Total	C	N	O	P	0	0
			140	70	14	49	7		

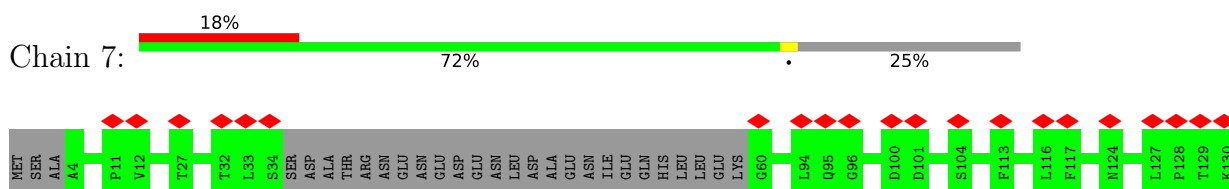
- Molecule 14 is a protein called DNA polymerase epsilon subunit B.

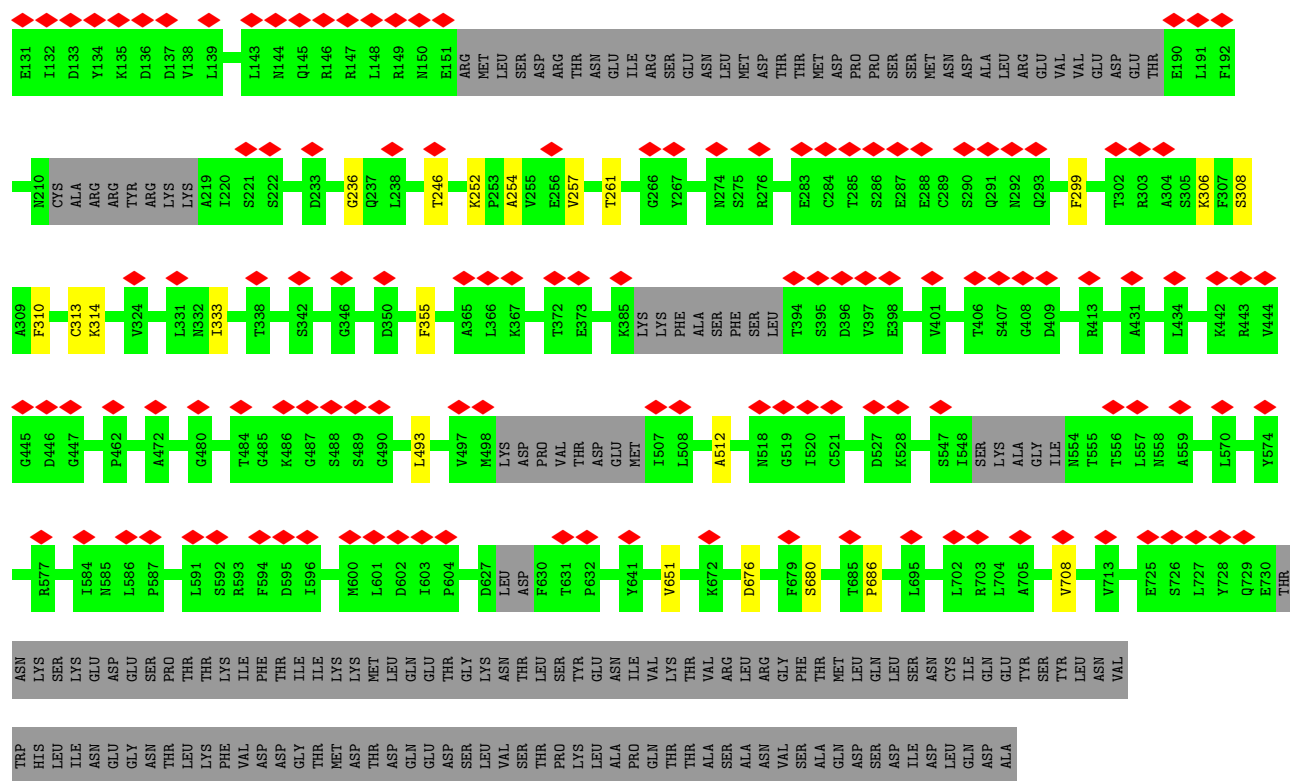
Mol	Chain	Residues	Atoms				AltConf	Trace
14	N	85	Total	C	N	O	0	0
			340	170	85	85		

Chain 6:

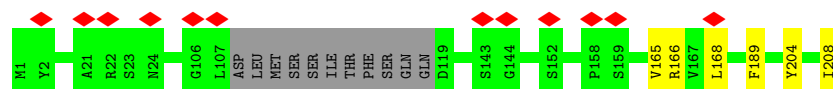
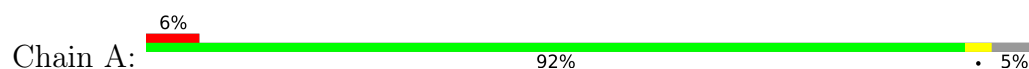


Chain 7:

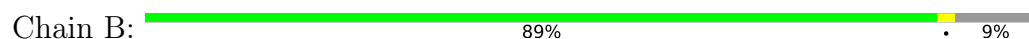




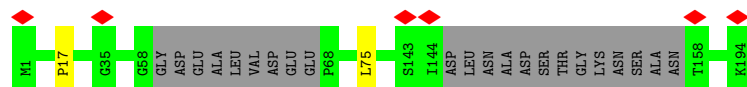
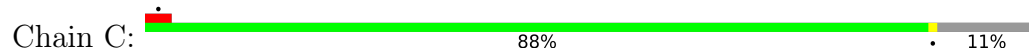
• Molecule 7: DNA replication complex GINS protein PSF1



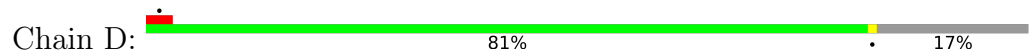
• Molecule 8: DNA replication complex GINS protein PSF2

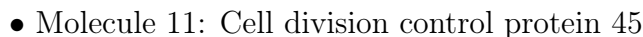


• Molecule 9: DNA replication complex GINS protein PSF3



• Molecule 10: DNA replication complex GINS protein SLD5





[illegible]

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	7939	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 ($e^-/\text{\AA}^2$)	53	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	81000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.416	Depositor
Minimum map value	-0.450	Depositor
Average map value	0.010	Depositor
Map value standard deviation	0.064	Depositor
Recommended contour level	0.45	Depositor
Map size ($\text{\AA}$)	466.39996, 466.39996, 466.39996	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.06, 1.06, 1.06	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	2	0.22	0/2461	0.43	2/3061 (0.1%)
2	3	0.16	0/2442	0.29	0/3042
3	4	0.12	0/2487	0.28	0/3100
4	5	0.22	0/2070	0.38	0/2563
5	6	0.09	0/2445	0.24	0/3044
6	7	0.17	0/2524	0.33	1/3141 (0.0%)
7	A	0.14	0/787	0.27	0/979
8	B	0.09	0/770	0.20	0/959
9	C	0.11	0/686	0.23	0/851
10	D	0.10	0/970	0.21	0/1206
11	E	0.11	0/2252	0.24	0/2805
12	F	0.08	0/1693	0.23	0/2111
12	G	0.08	0/1685	0.23	0/2101
12	H	0.08	0/1697	0.22	0/2116
13	I	0.19	0/153	0.51	0/234
14	N	0.11	0/339	0.34	0/422
All	All	0.14	0/25461	0.30	3/31735 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	7	686	PRO	N-CA-C	-7.00	103.80	113.53
1	2	627	GLN	N-CA-C	-6.00	101.28	113.29
1	2	622	GLU	N-CA-C	-5.74	105.02	111.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	2	2469	0	648	7	0
2	3	2448	0	642	9	0
3	4	2492	0	648	8	0
4	5	2084	0	528	11	0
5	6	2452	0	647	4	0
6	7	2532	0	662	10	0
7	A	789	0	213	3	0
8	B	772	0	189	2	0
9	C	689	0	175	1	0
10	D	973	0	240	2	0
11	E	2257	0	603	8	0
12	F	1696	0	444	9	0
12	G	1688	0	442	7	0
12	H	1700	0	445	14	0
13	I	140	0	85	2	0
14	N	340	0	94	1	0
All	All	25521	0	6705	97	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.

All (97) 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)
13:I:6:DT:H2"	13:I:7:DT:H5"	1.74	0.68
11:E:64:TYR:N	11:E:623:ASP:O	2.27	0.67
4:5:175:ARG:N	4:5:251:ILE:O	2.22	0.66
6:7:236:GLY:N	6:7:355:PHE:O	2.31	0.63
3:4:604:TYR:O	3:4:617:GLU:N	2.29	0.63
10:D:224:TRP:O	10:D:280:GLU:N	2.31	0.63
3:4:178:ARG:O	3:4:187:ILE:N	2.34	0.61
6:7:246:THR:N	6:7:314:LYS:O	2.33	0.61
8:B:117:TRP:O	8:B:175:LEU:N	2.28	0.61
1:2:315:SER:N	1:2:430:TYR:O	2.30	0.61
11:E:579:TYR:N	11:E:632:ILE:O	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:E:323:ASP:N	11:E:406:ARG:O	2.34	0.59
12:H:840:THR:O	12:H:845:GLY:N	2.36	0.59
6:7:254:ALA:N	6:7:308:SER:O	2.34	0.58
11:E:315:THR:O	11:E:414:GLY:N	2.36	0.58
2:3:406:LEU:O	2:3:547:PHE:N	2.35	0.58
2:3:317:PHE:N	4:5:174:SER:O	2.25	0.58
12:H:601:ASN:N	12:H:605:VAL:O	2.37	0.58
11:E:33:CYS:N	11:E:61:ILE:O	2.31	0.56
12:H:577:ARG:O	12:H:593:SER:N	2.36	0.56
12:G:640:LEU:N	12:G:652:LYS:O	2.39	0.55
4:5:173:SER:N	4:5:253:GLN:O	2.39	0.55
12:H:597:PHE:N	12:H:610:GLU:O	2.40	0.55
12:H:743:TRP:N	12:H:755:ILE:O	2.38	0.55
6:7:252:LYS:O	6:7:310:PHE:N	2.37	0.55
5:6:301:ARG:O	5:6:356:TRP:N	2.38	0.55
12:H:638:TYR:N	12:H:655:CYS:O	2.39	0.55
12:F:479:PHE:N	12:F:579:THR:O	2.39	0.54
7:A:165:VAL:O	7:A:189:PHE:N	2.41	0.53
4:5:53:ASN:O	4:5:58:ASN:N	2.39	0.53
12:H:638:TYR:O	12:H:655:CYS:N	2.42	0.53
12:G:597:PHE:N	12:G:610:GLU:O	2.34	0.53
3:4:565:LEU:O	3:4:706:TYR:N	2.37	0.53
5:6:398:THR:O	5:6:457:CYS:N	2.42	0.53
1:2:433:ASN:O	1:2:448:ALA:N	2.42	0.52
5:6:636:CYS:N	5:6:677:SER:O	2.41	0.52
4:5:180:SER:O	4:5:244:ILE:N	2.40	0.52
12:F:505:VAL:O	12:F:512:SER:N	2.43	0.52
11:E:328:LEU:O	11:E:332:SER:N	2.38	0.51
12:F:741:HIS:N	12:F:757:VAL:O	2.45	0.50
2:3:433:THR:O	2:3:437:SER:N	2.45	0.49
4:5:179:LEU:N	4:5:192:ILE:O	2.46	0.49
1:2:604:CYS:N	1:2:645:SER:O	2.45	0.49
3:4:593:GLY:N	3:4:632:ASP:O	2.45	0.49
12:G:638:TYR:O	12:G:655:CYS:N	2.46	0.48
4:5:157:SER:N	4:5:298:TYR:O	2.46	0.48
12:F:588:VAL:N	12:F:600:PHE:O	2.36	0.48
9:C:17:PRO:O	9:C:75:LEU:N	2.45	0.48
12:G:629:HIS:O	12:G:637:SER:N	2.47	0.48
2:3:199:SER:N	2:3:212:ARG:O	2.43	0.48
12:G:501:TYR:N	12:G:516:SER:O	2.40	0.48
4:5:177:THR:N	4:5:249:LYS:O	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:N:34:LYS:H	14:N:76:GLY:HA3	1.77	0.47
2:3:177:ASN:N	2:3:298:PHE:O	2.47	0.47
2:3:166:LEU:N	2:3:181:SER:O	2.47	0.47
3:4:600:GLY:O	3:4:620:ALA:N	2.45	0.47
5:6:307:ALA:O	5:6:321:VAL:N	2.35	0.47
12:H:746:ALA:O	12:H:753:ASN:N	2.47	0.46
10:D:226:LYS:O	10:D:279:TYR:N	2.47	0.46
12:H:685:LYS:N	12:H:698:PHE:O	2.47	0.46
1:2:407:GLU:O	1:2:451:ILE:N	2.48	0.46
12:F:597:PHE:O	12:F:610:GLU:N	2.45	0.46
6:7:651:VAL:O	6:7:708:VAL:N	2.40	0.45
2:3:301:LEU:N	2:3:318:LYS:O	2.48	0.45
7:A:168:LEU:N	7:A:204:TYR:O	2.49	0.45
3:4:437:GLY:HA2	3:4:464:VAL:H	1.81	0.45
12:G:577:ARG:O	12:G:593:SER:N	2.42	0.45
1:2:600:ASP:C	1:2:602:GLY:H	2.25	0.45
2:3:434:GLY:N	2:3:473:ASP:O	2.46	0.44
8:B:126:LEU:O	8:B:130:ALA:N	2.51	0.44
12:H:627:SER:O	12:H:639:SER:N	2.44	0.44
3:4:434:GLU:N	3:4:467:LYS:O	2.50	0.43
12:H:492:ARG:N	12:H:504:THR:O	2.47	0.43
1:2:653:ASN:N	1:2:666:ASN:O	2.47	0.43
11:E:100:PRO:O	11:E:104:VAL:N	2.48	0.43
12:F:546:PHE:N	12:F:555:GLN:O	2.46	0.42
4:5:182:MET:N	4:5:242:ILE:O	2.46	0.42
6:7:313:CYS:O	6:7:333:ILE:N	2.52	0.42
12:G:746:ALA:O	12:G:753:ASN:N	2.51	0.42
12:F:704:LEU:N	12:F:722:LEU:O	2.52	0.42
12:H:741:HIS:N	12:H:757:VAL:O	2.53	0.41
12:F:546:PHE:O	12:F:555:GLN:N	2.45	0.41
12:H:500:GLY:HA2	12:H:518:PHE:H	1.86	0.41
2:3:434:GLY:HA3	2:3:474:GLU:O	2.20	0.41
4:5:439:THR:O	4:5:480:ASP:N	2.38	0.41
3:4:204:LYS:N	3:4:221:ASP:O	2.54	0.41
11:E:637:LEU:O	11:E:641:LEU:N	2.48	0.41
12:F:599:SER:O	12:F:607:PHE:N	2.54	0.41
4:5:613:ARG:O	4:5:672:ALA:N	2.31	0.41
6:7:257:VAL:O	6:7:306:LYS:N	2.53	0.41
6:7:676:ASP:O	6:7:680:SER:N	2.54	0.41
1:2:305:SER:O	1:2:309:LEU:N	2.51	0.41
12:H:629:HIS:N	12:H:637:SER:O	2.43	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:166:ARG:N	7:A:208:ILE:CA	2.84	0.40
13:I:5:DT:H6	13:I:5:DT:H2'	1.74	0.40
6:7:261:THR:O	6:7:299:PHE:N	2.45	0.40
6:7:493:LEU:O	6:7:512:ALA:N	2.54	0.40

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	2	601/868 (69%)	590 (98%)	11 (2%)	0	100	100
2	3	600/971 (62%)	591 (98%)	9 (2%)	0	100	100
3	4	613/933 (66%)	605 (99%)	8 (1%)	0	100	100
4	5	493/775 (64%)	484 (98%)	9 (2%)	0	100	100
5	6	599/1017 (59%)	587 (98%)	12 (2%)	0	100	100
6	7	617/845 (73%)	611 (99%)	6 (1%)	0	100	100
7	A	193/208 (93%)	191 (99%)	2 (1%)	0	100	100
8	B	189/213 (89%)	184 (97%)	5 (3%)	0	100	100
9	C	166/194 (86%)	164 (99%)	2 (1%)	0	100	100
10	D	237/294 (81%)	235 (99%)	2 (1%)	0	100	100
11	E	554/650 (85%)	549 (99%)	5 (1%)	0	100	100
12	F	418/927 (45%)	409 (98%)	9 (2%)	0	100	100
12	G	416/927 (45%)	410 (99%)	6 (1%)	0	100	100
12	H	419/927 (45%)	410 (98%)	9 (2%)	0	100	100
14	N	83/689 (12%)	80 (96%)	3 (4%)	0	100	100
All	All	6198/10438 (59%)	6100 (98%)	98 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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.

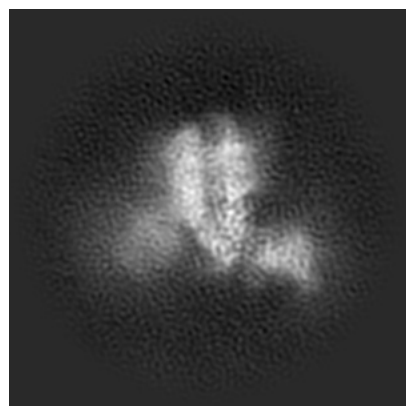
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-37345. These allow visual inspection of the internal detail of the map and identification of artifacts.

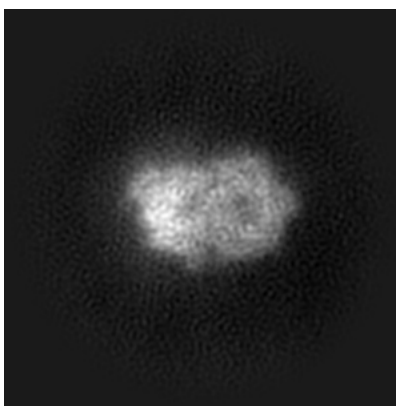
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

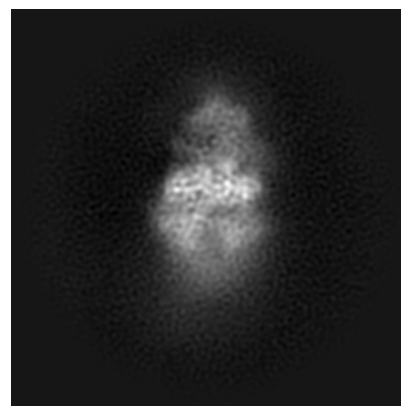
6.1.1 Primary map



X

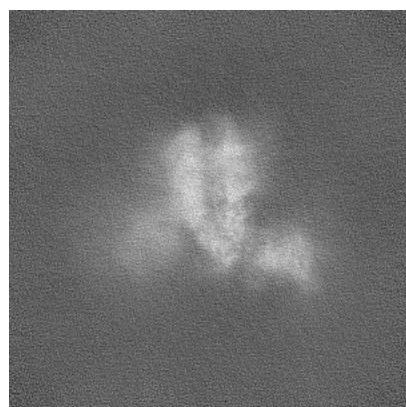


Y

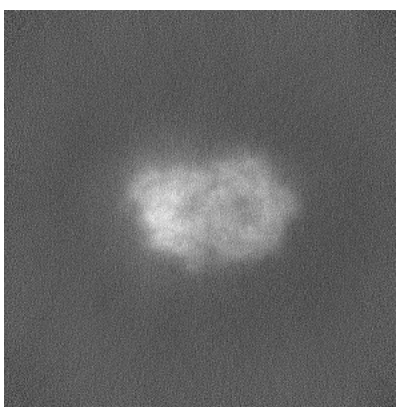


Z

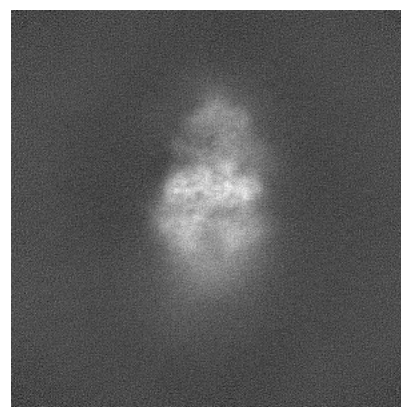
6.1.2 Raw map



X



Y

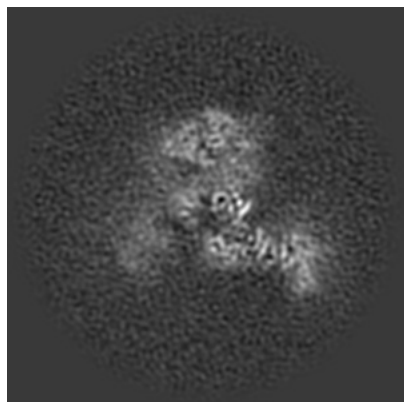


Z

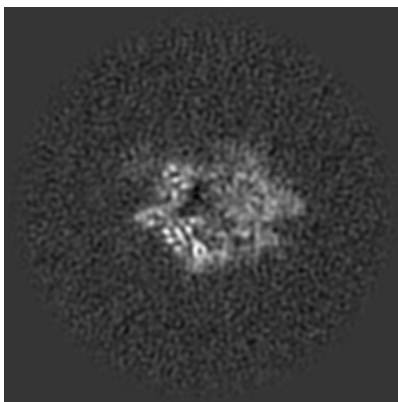
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

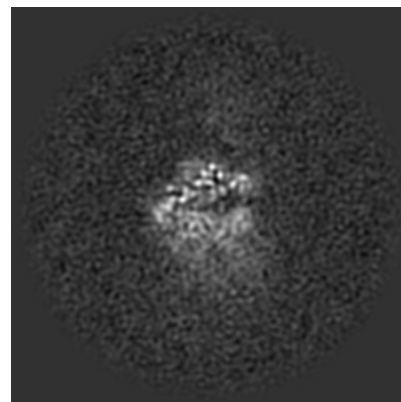
6.2.1 Primary map



X Index: 220

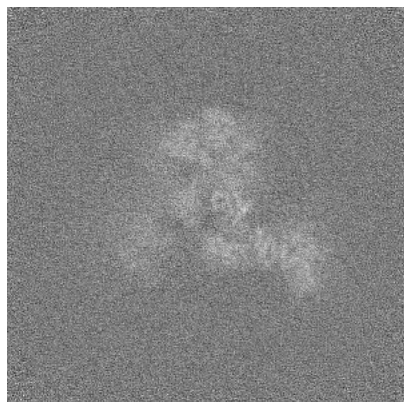


Y Index: 220

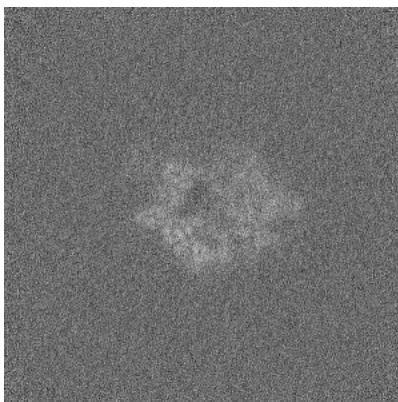


Z Index: 220

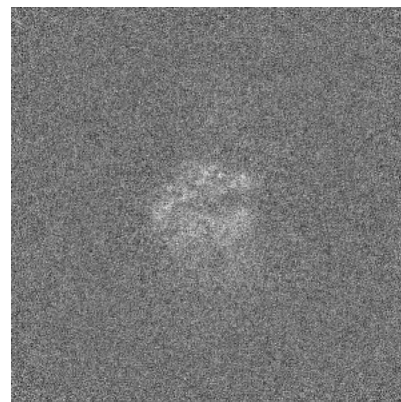
6.2.2 Raw map



X Index: 220



Y Index: 220

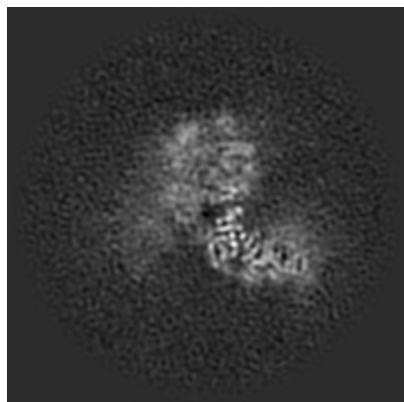


Z Index: 220

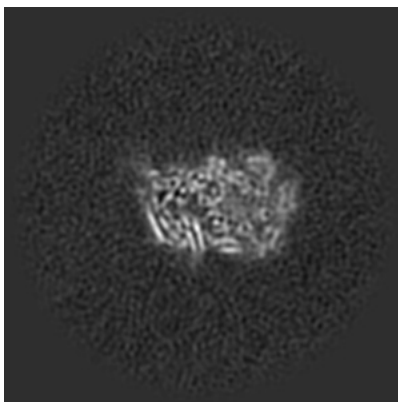
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

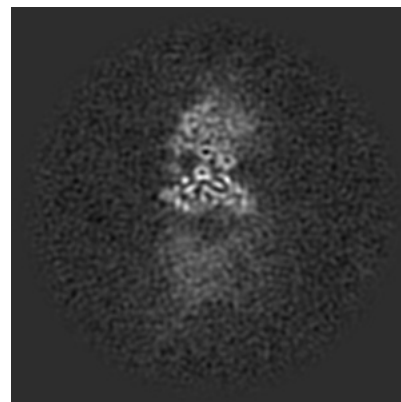
6.3.1 Primary map



X Index: 238

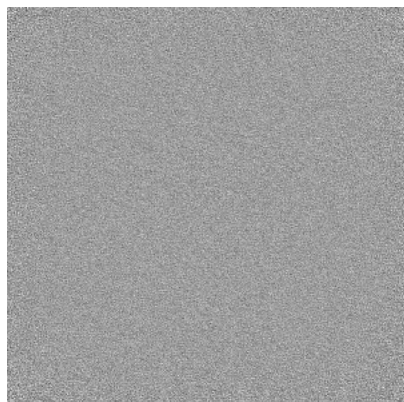


Y Index: 242

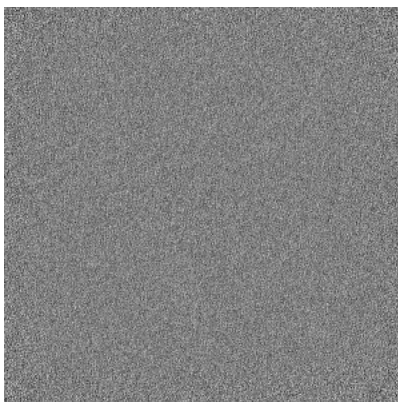


Z Index: 179

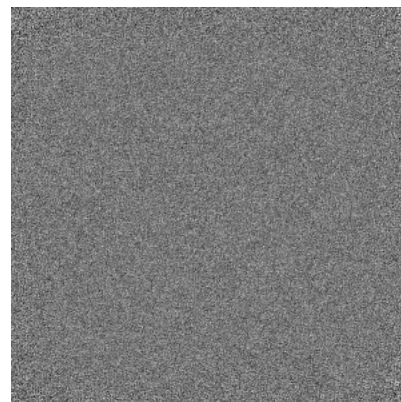
6.3.2 Raw map



X Index: 0



Y Index: 0

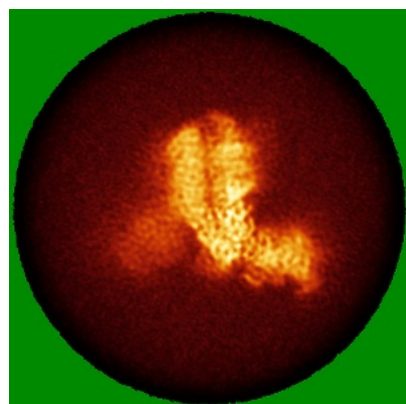


Z Index: 0

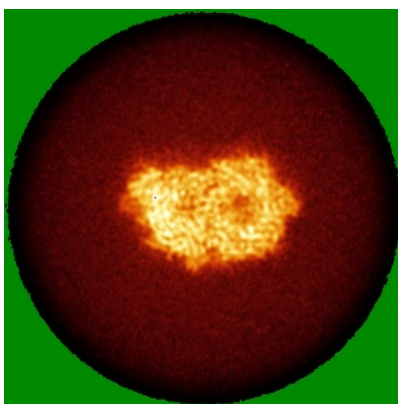
The images above show the largest variance slices of the map in three orthogonal directions.

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

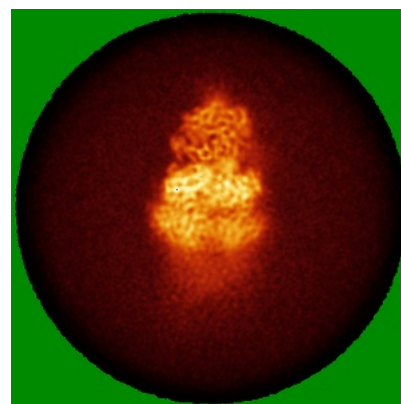
6.4.1 Primary map



X

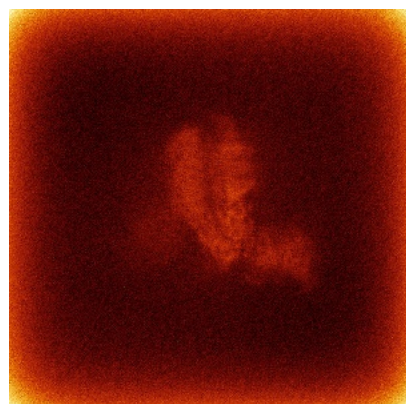


Y

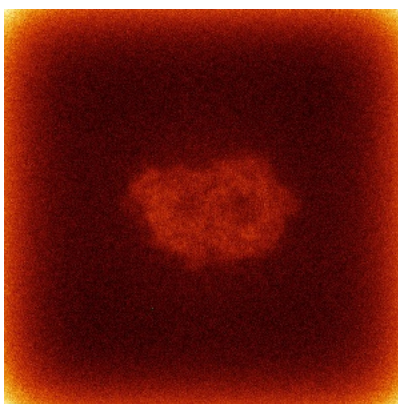


Z

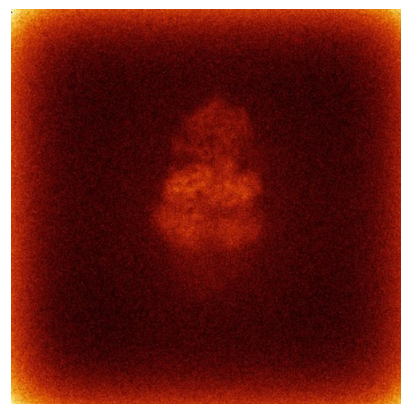
6.4.2 Raw map



X



Y

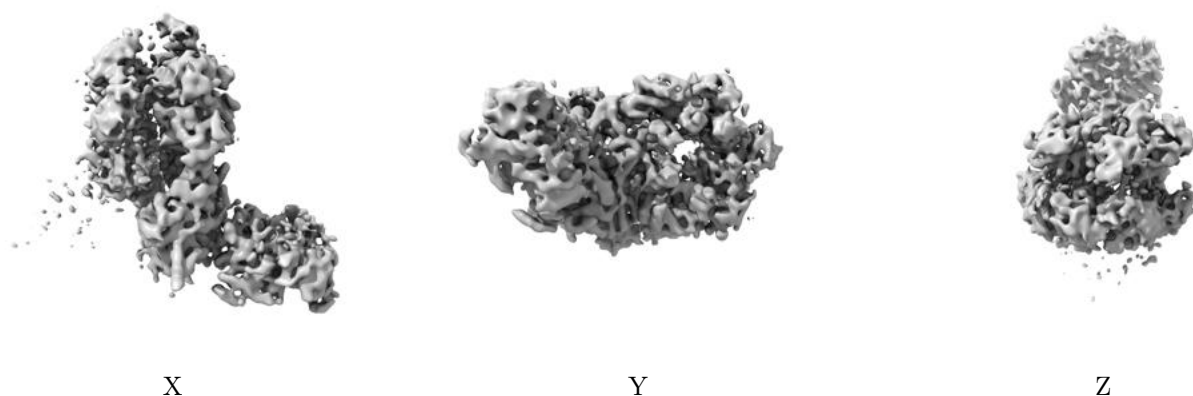


Z

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.

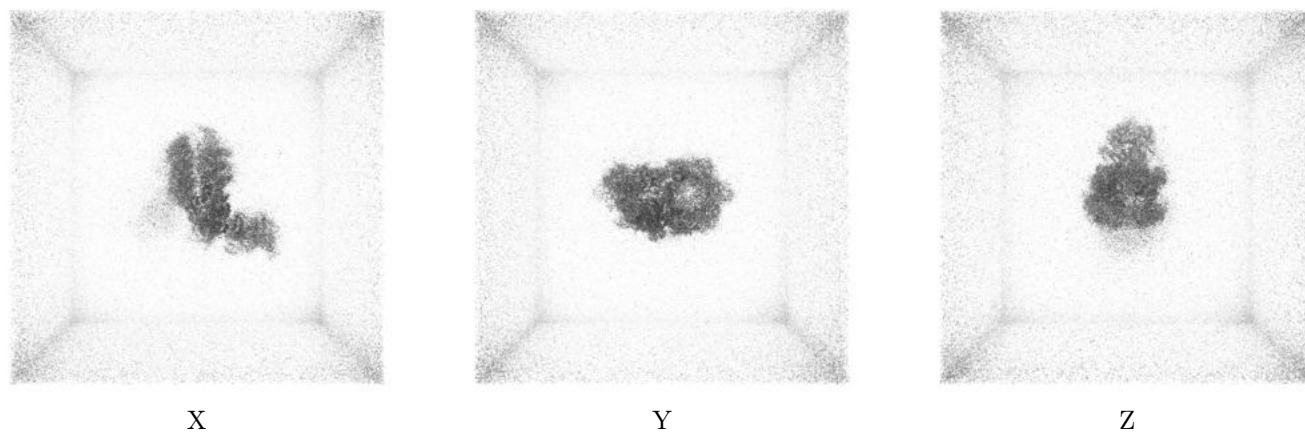
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.45. 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.

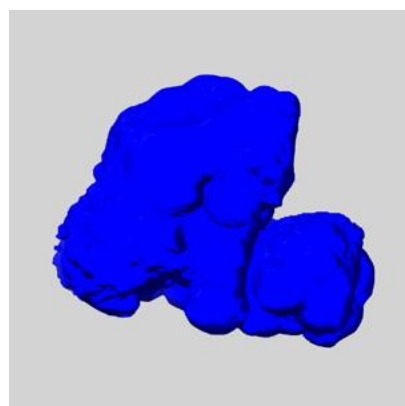
6.6 Mask visualisation [i](#)

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

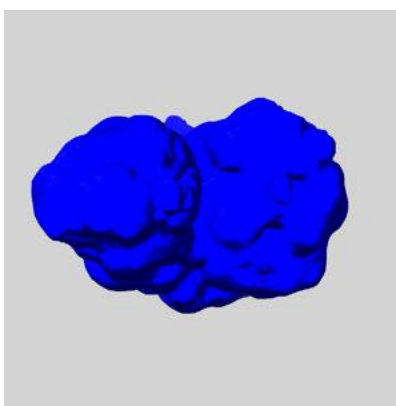
A mask typically either:

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

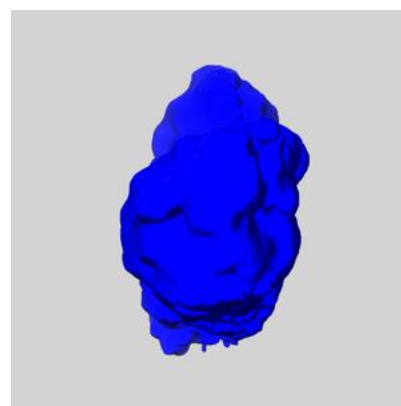
6.6.1 emd_37345_msk_1.map [i](#)



X



Y

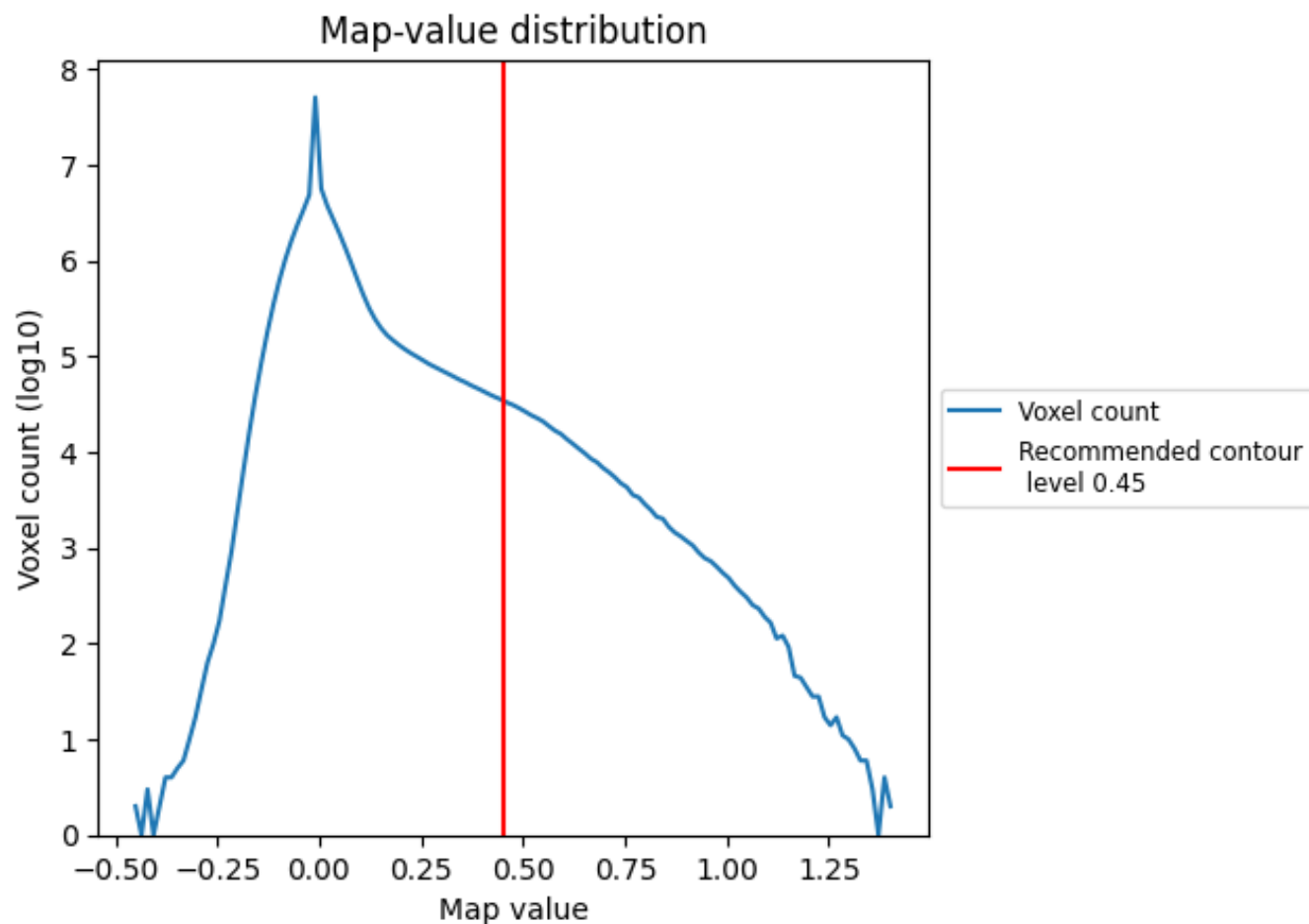


Z

7 Map analysis [i](#)

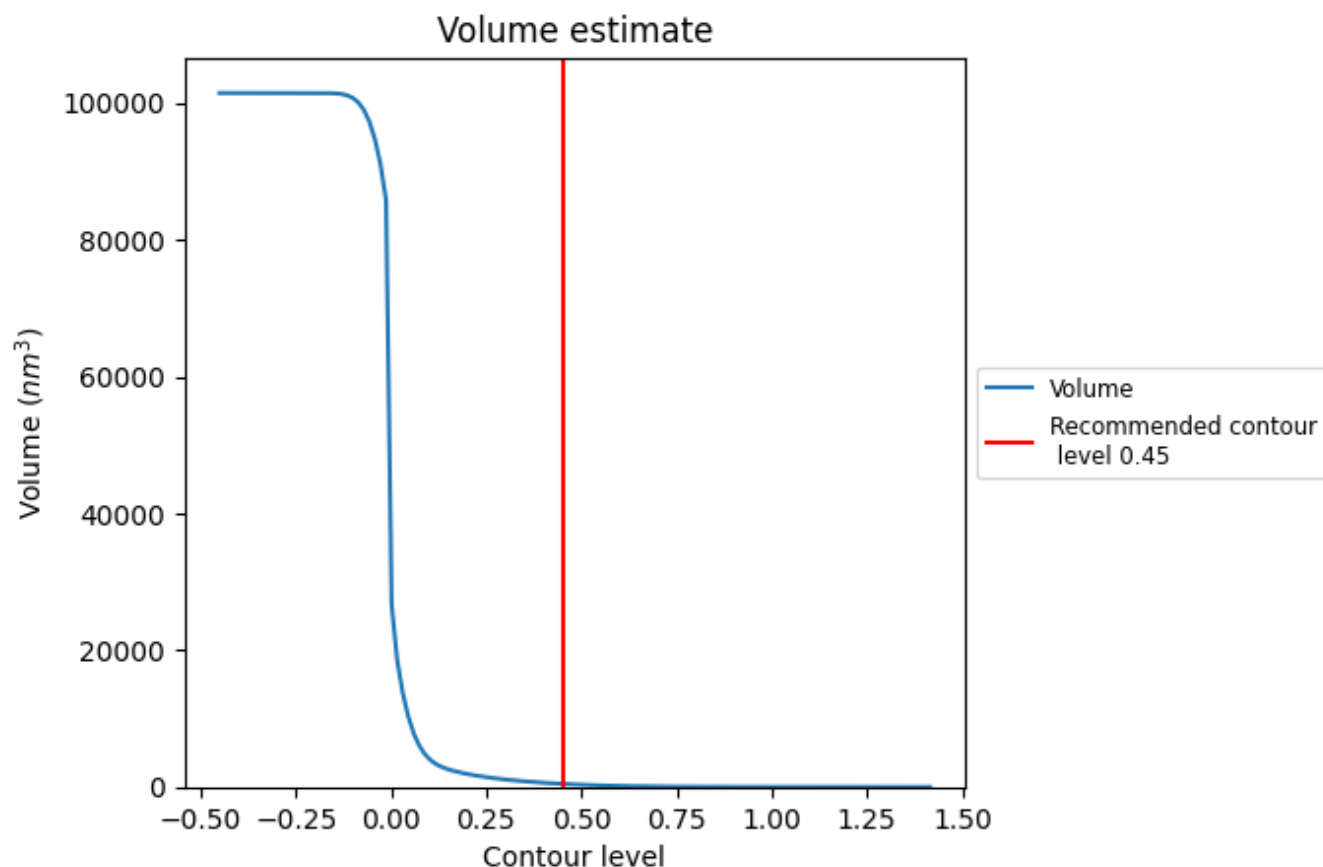
This section contains the results of statistical analysis of the map.

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.

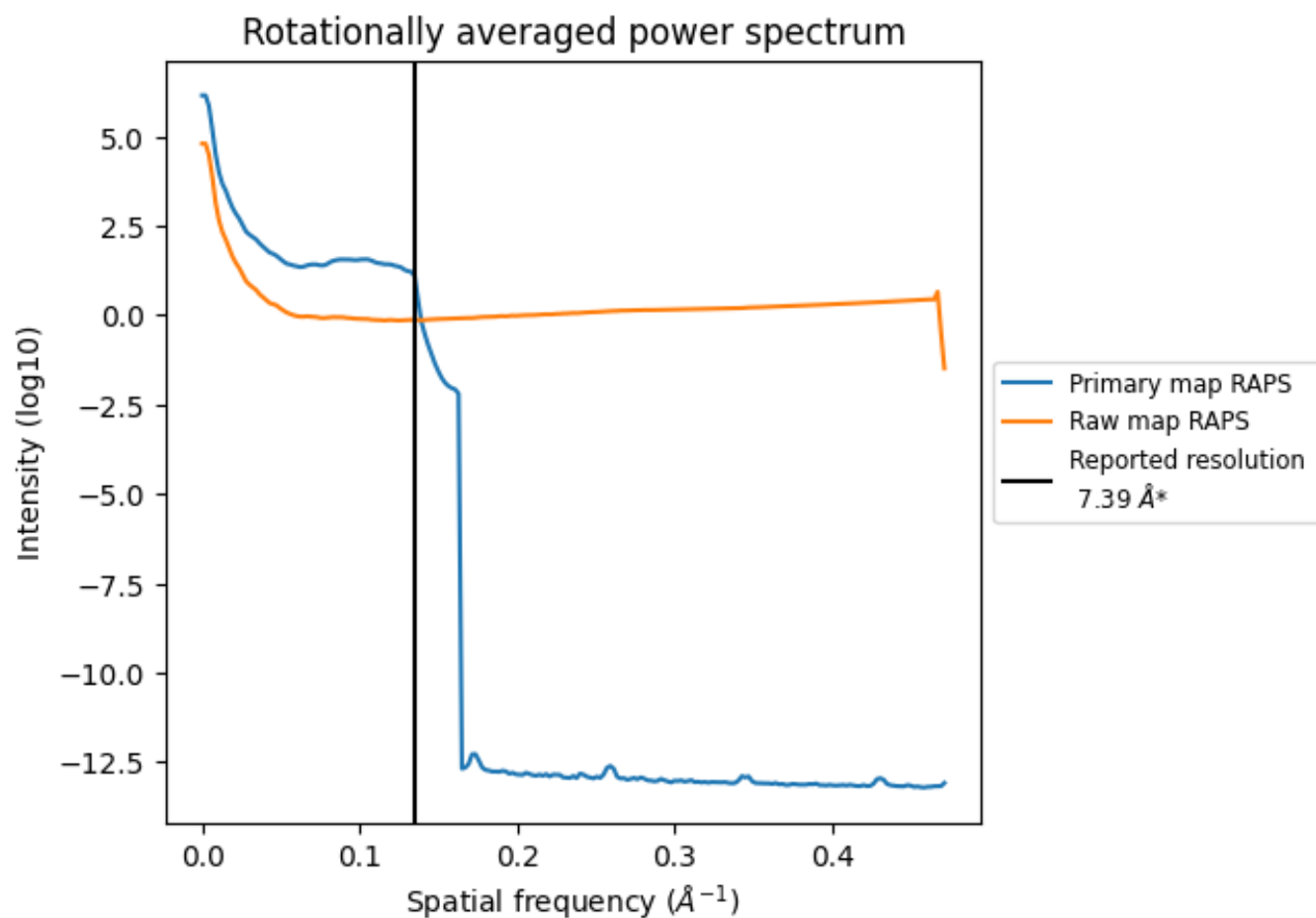
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 457 nm^3 ; this corresponds to an approximate mass of 412 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

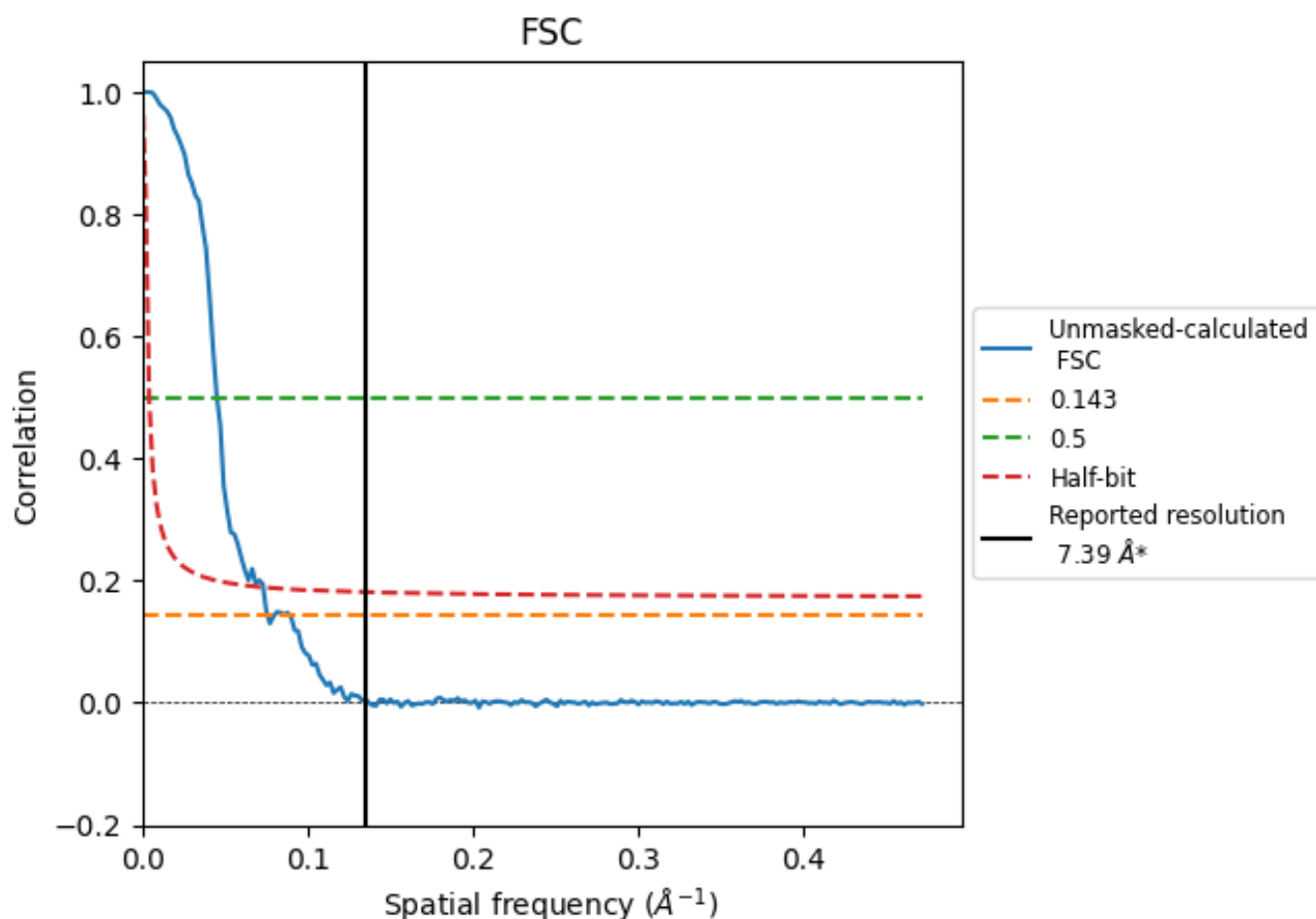


*Reported resolution corresponds to spatial frequency of 0.135 Å⁻¹

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 $0.135 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

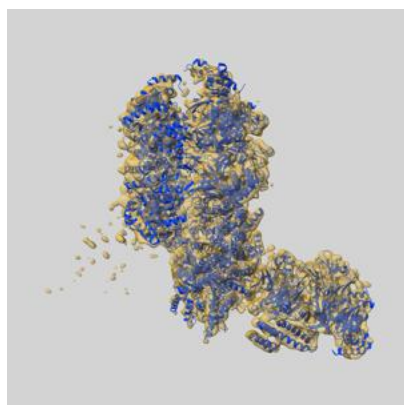
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	7.39	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	13.19	22.12	13.66

*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 13.19 differs from the reported value 7.39 by more than 10 %

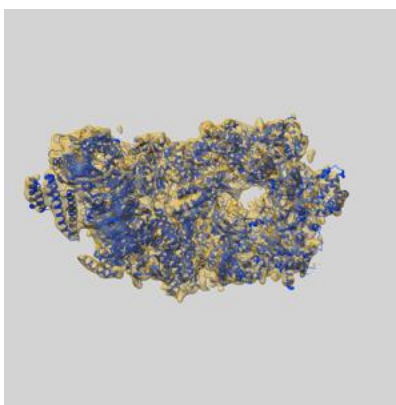
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-37345 and PDB model 8W7S. Per-residue inclusion information can be found in section 3 on page 7.

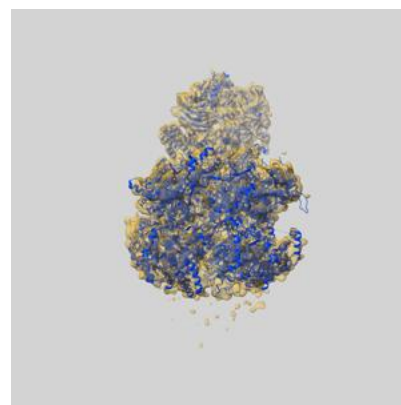
9.1 Map-model overlay [i](#)



X



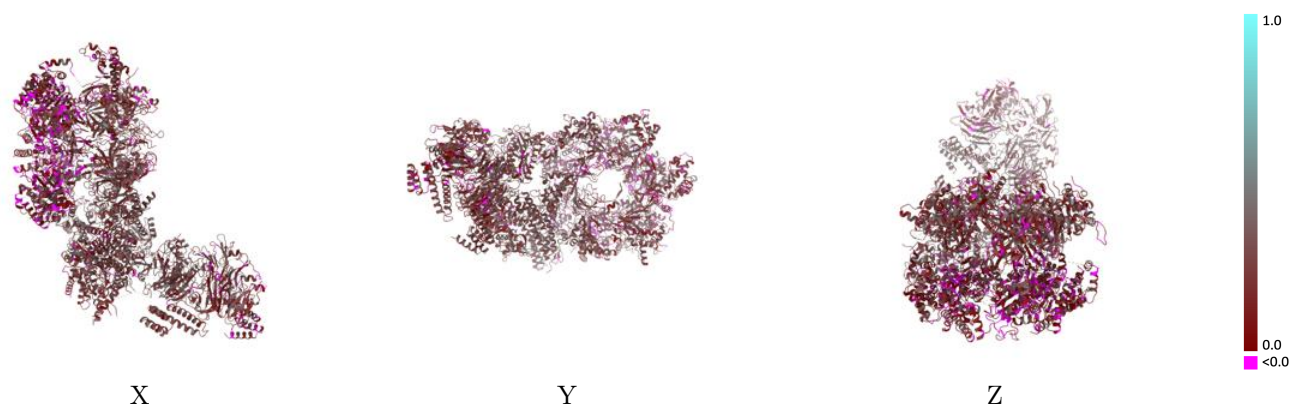
Y



Z

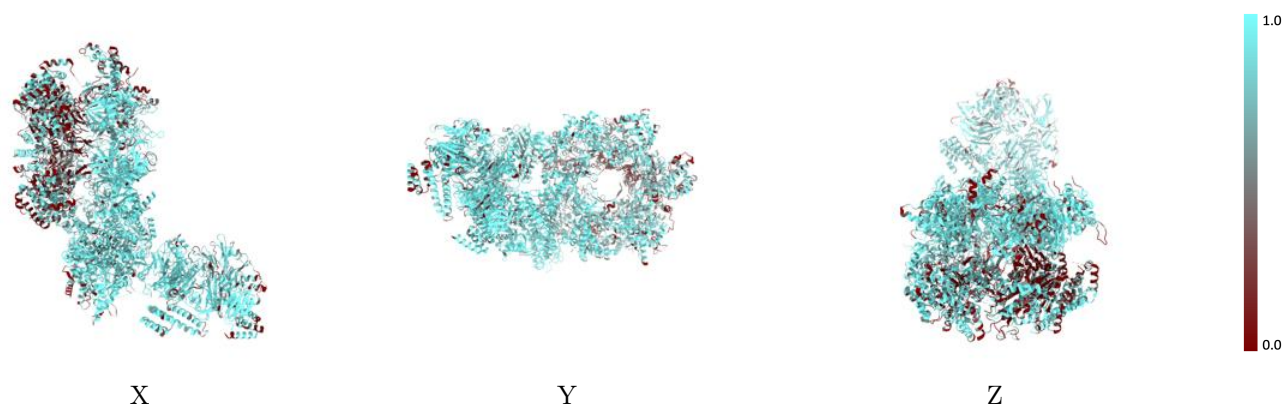
The images above show the 3D surface view of the map at the recommended contour level 0.45 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



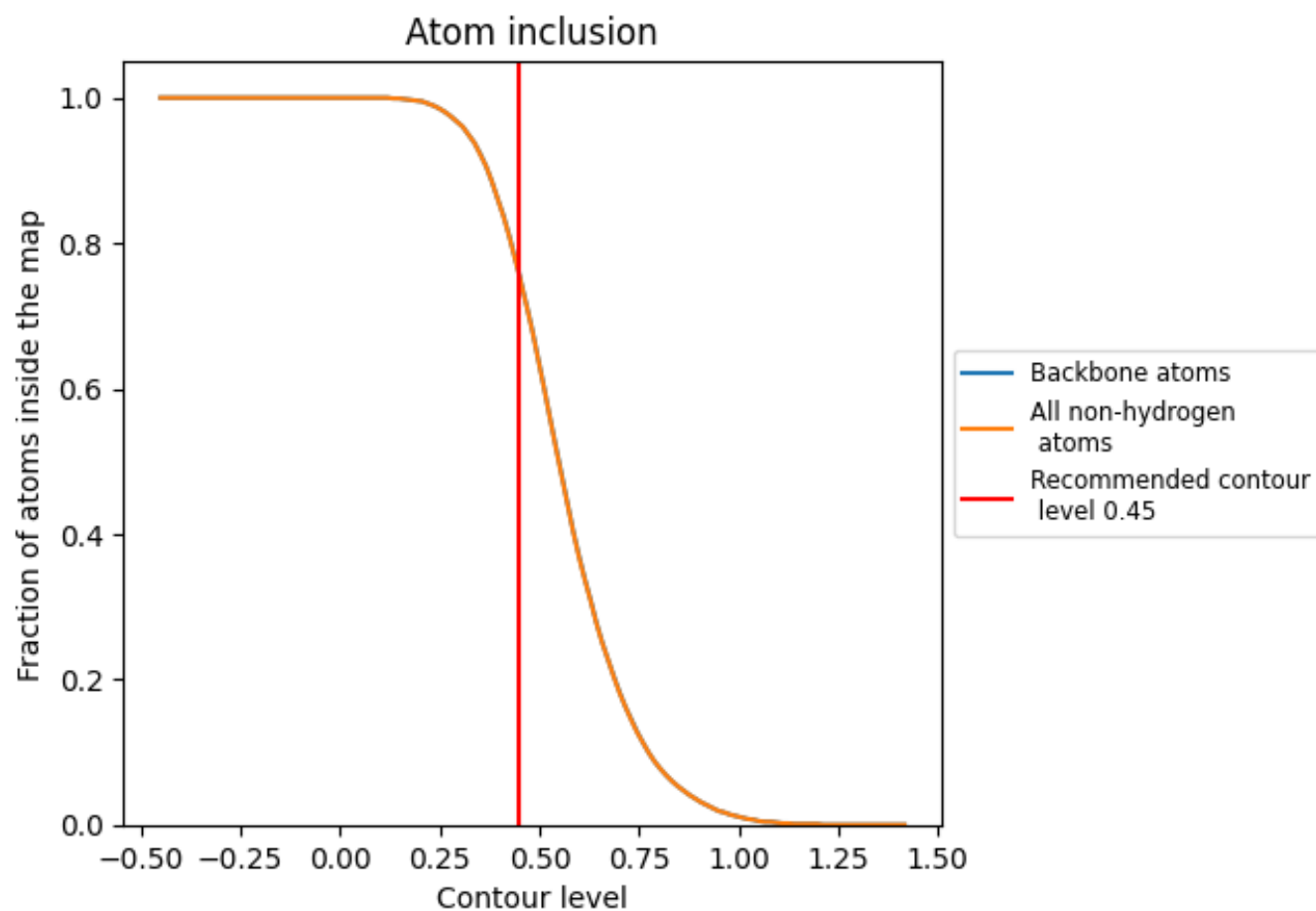
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.45).

9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 76% 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 (0.45) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7550	 0.2120
2	 0.6360	 0.1580
3	 0.6960	 0.1670
4	 0.5920	 0.1500
5	 0.8200	 0.2380
6	 0.5970	 0.1680
7	 0.6940	 0.1900
A	 0.9140	 0.2780
B	 0.9790	 0.3090
C	 0.9420	 0.2730
D	 0.9200	 0.2780
E	 0.8850	 0.2490
F	 0.9240	 0.2910
G	 0.7690	 0.2210
H	 0.8380	 0.2410
I	 0.5790	 0.1780
N	 0.5820	 0.2270

