



## Full wwPDB EM Validation Report ⓘ

May 11, 2026 – 11:07 PM EDT

PDB ID : 9DX0 / pdb\_00009dx0  
EMDB ID : EMD-47276  
Title : Human GATOR2 complex - apo state  
Authors : Wranik, M.; Rogala, K.B.  
Deposited on : 2024-10-10  
Resolution : 3.47 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

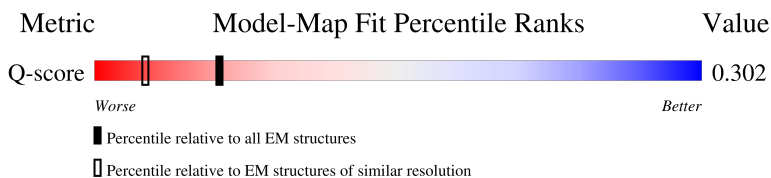
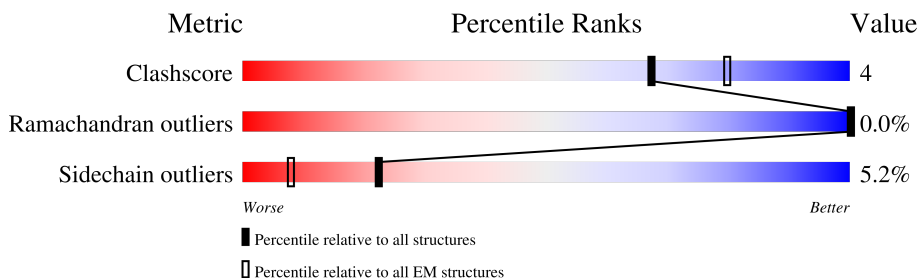
EMDB validation analysis : 0.0.1.dev132  
MolProbity : 4-5-2 with Phenix2.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)  
MapQ : 1.9.13  
Ideal geometry (proteins)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

# 1 Overall quality at a glance

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

The reported resolution of this entry is 3.47 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	13733 ( 2.97 - 3.97 )

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion  $< 40\%$ ). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1113	<div> <div>10%</div> <div>88%</div> <div>10%</div> <div>..</div> </div>
1	B	1113	<div> <div>18%</div> <div>86%</div> <div>10%</div> <div>..</div> </div>
2	C	905	<div> <div>39%</div> <div>89%</div> <div>10%</div> <div>.</div> </div>
3	D	633	<div> <div>6%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 27053 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 GATOR2 complex protein MIOS.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1103	Total	C	N	O	S	0	0
			8290	5252	1465	1512	61		
1	B	1086	Total	C	N	O	S	0	0
			7352	4595	1340	1374	43		

- Molecule 2 is a protein called GATOR2 complex protein WDR24.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	905	Total	C	N	O	S	0	0
			6746	4241	1222	1223	60		

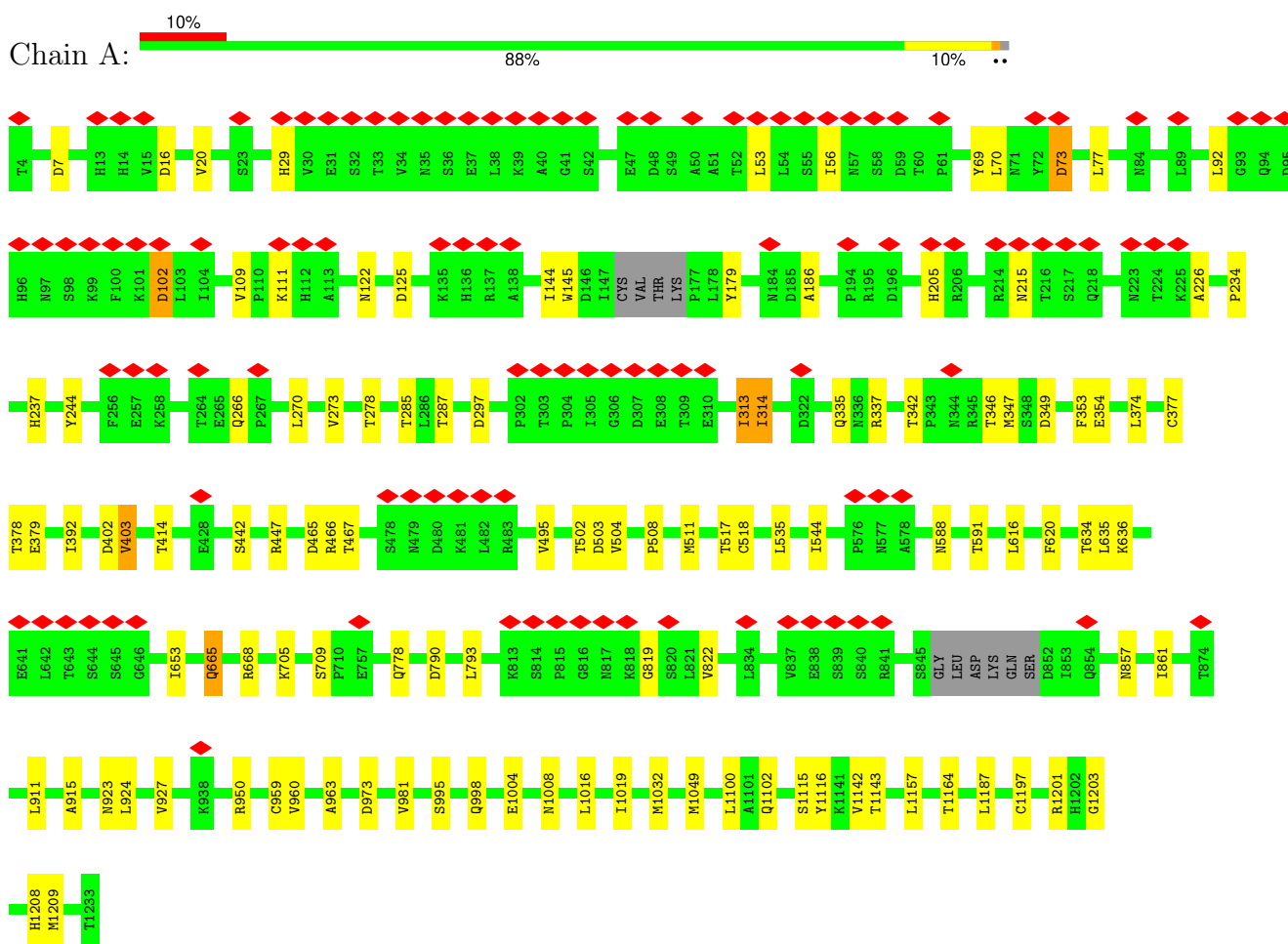
- Molecule 3 is a protein called Protein SEC13 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	633	Total	C	N	O	S	0	0
			4665	2935	831	863	36		

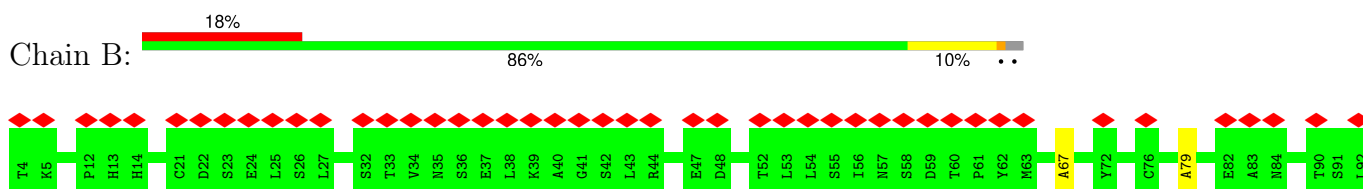
### 3 Residue-property plots [i](#)

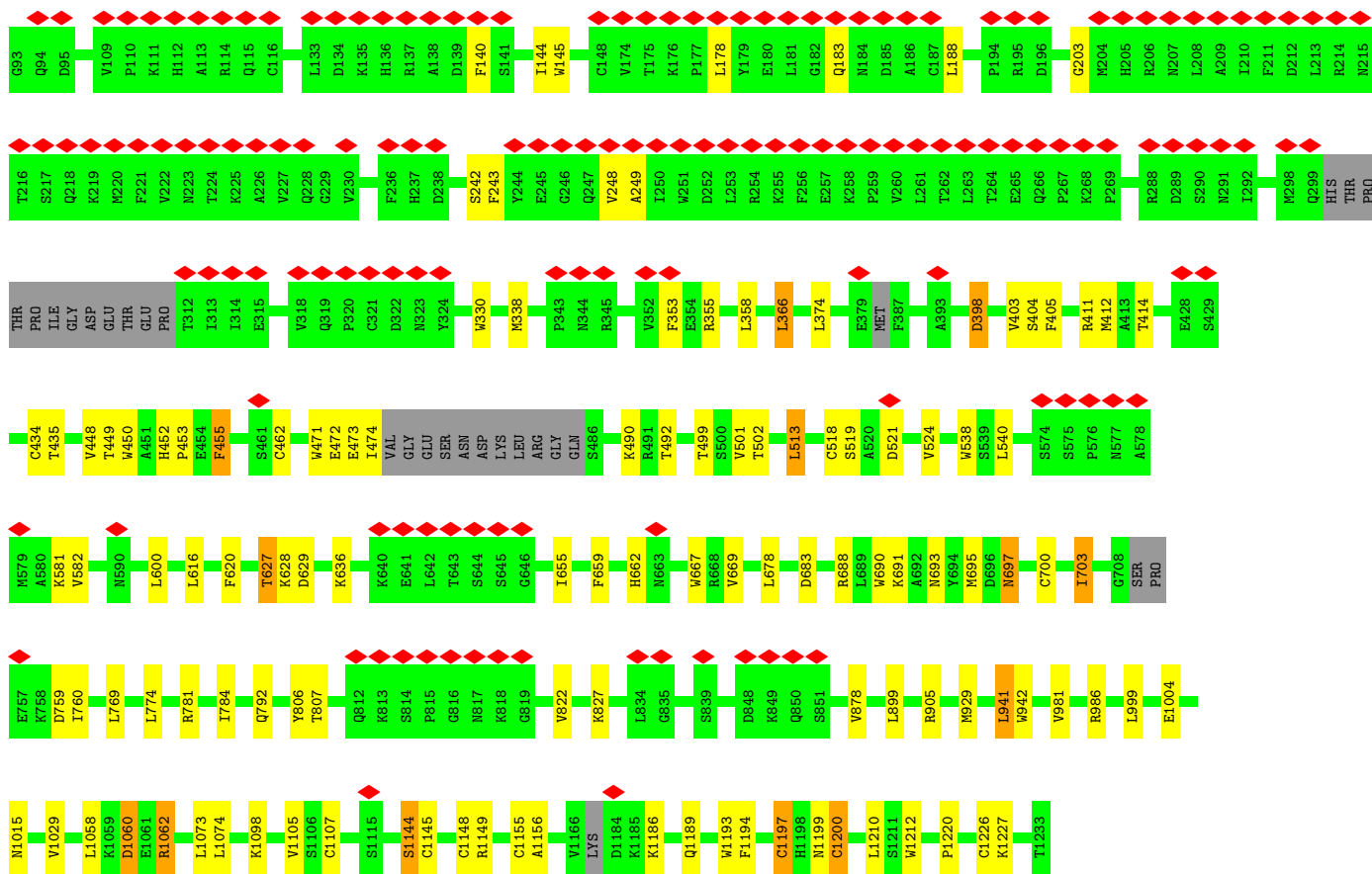
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: GATOR2 complex protein MIOS

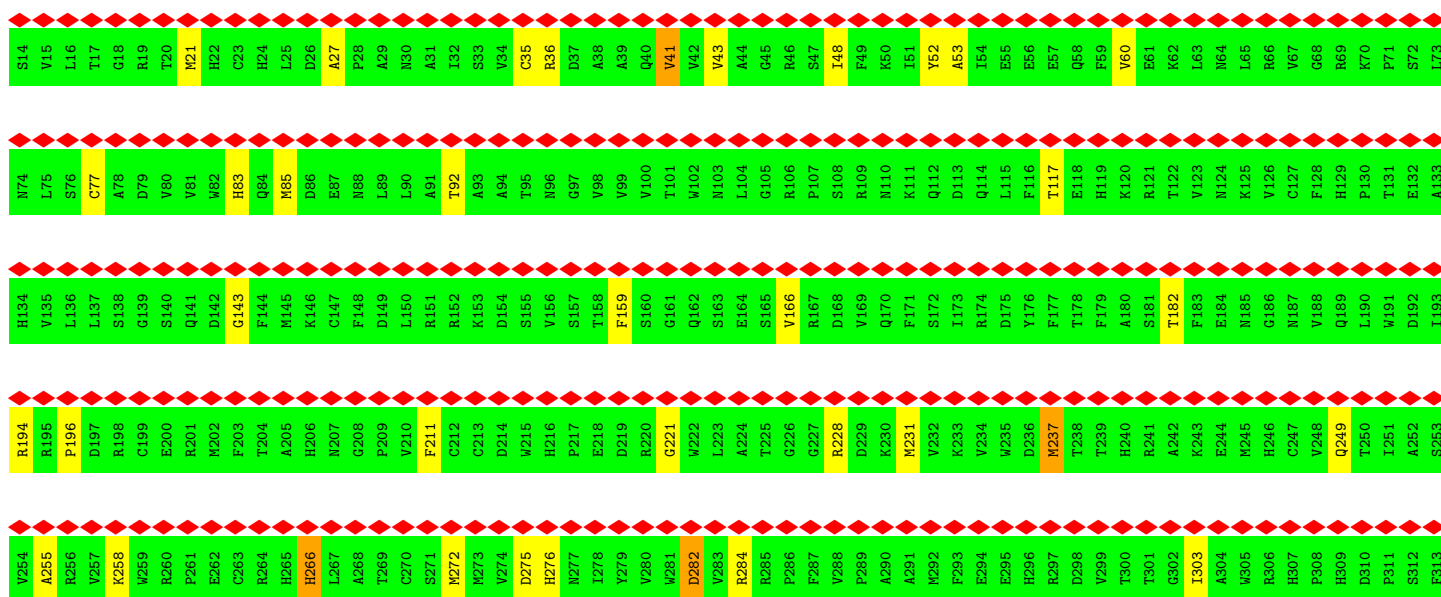
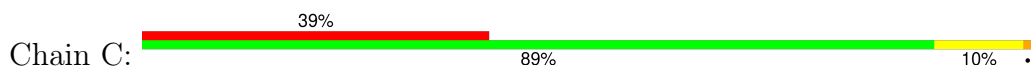


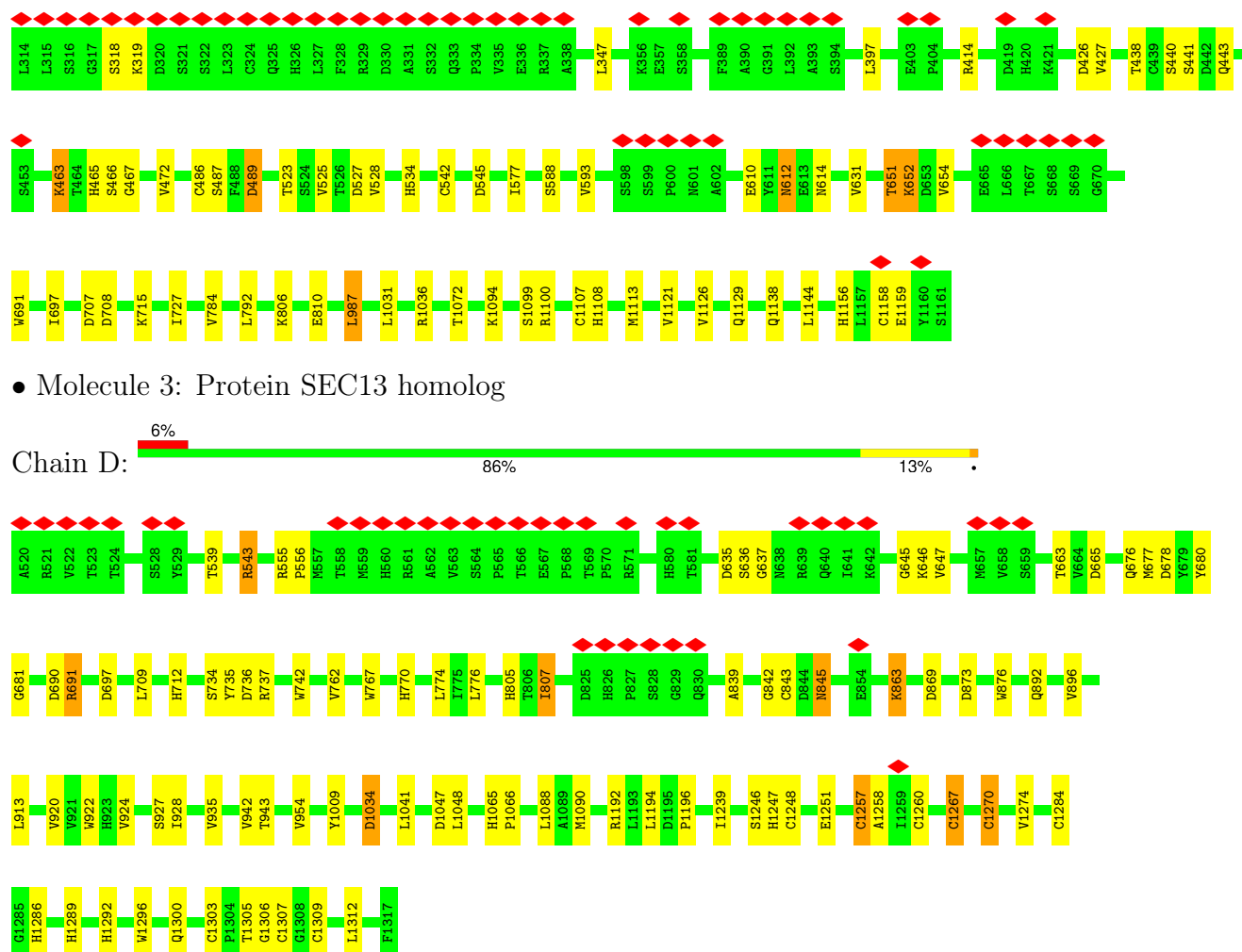
#### • Molecule 1: GATOR2 complex protein MIOS





• Molecule 2: GATOR2 complex protein WDR24





• Molecule 3: Protein SEC13 homolog

## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	1148926	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 ( $e^-/\text{\AA}^2$ )	64.64	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2100	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.432	Depositor
Minimum map value	-0.154	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.011	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	417.456, 417.456, 417.456	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8697, 0.8697, 0.8697	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	A	0.17	0/8492	0.27	0/11586
1	B	0.16	0/7488	0.27	0/10252
2	C	0.18	0/6921	0.27	0/9440
3	D	0.21	0/4780	0.30	0/6516
All	All	0.18	0/27681	0.28	0/37794

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	8290	0	7731	56	0
1	B	7352	0	6068	60	0
2	C	6746	0	6204	42	0
3	D	4665	0	4280	42	0
All	All	27053	0	24283	195	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 4.

All (195) 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)
3:D:1307:CYS:SG	3:D:1309:CYS:N	2.58	0.75
2:C:1156:HIS:CE1	2:C:1158:CYS:HB2	2.22	0.74
3:D:1270:CYS:SG	3:D:1289:HIS:ND1	2.58	0.74
1:A:278:THR:HG21	1:A:335:GLN:HA	1.70	0.74
3:D:1267:CYS:SG	3:D:1270:CYS:N	2.58	0.71
1:B:1197:CYS:SG	1:B:1200:CYS:N	2.61	0.70
2:C:612:ASN:ND2	2:C:614:ASN:OD1	2.26	0.69
3:D:1270:CYS:SG	3:D:1292:HIS:ND1	2.66	0.68
1:B:1145:CYS:HB3	1:B:1148:CYS:SG	2.35	0.67
3:D:1286:HIS:HE1	3:D:1307:CYS:HB3	1.59	0.67
1:A:73:ASP:OD1	1:A:73:ASP:N	2.27	0.67
1:A:186:ALA:HB3	1:A:205:HIS:HB2	1.77	0.67
3:D:646:LYS:NZ	3:D:665:ASP:OD1	2.29	0.65
3:D:892:GLN:HA	3:D:920:VAL:HG23	1.78	0.64
1:A:53:LEU:HD11	1:A:56:ILE:HD11	1.79	0.64
1:A:950:ARG:NH2	1:A:959:CYS:SG	2.71	0.63
1:A:402:ASP:OD1	1:A:403:VAL:N	2.32	0.62
1:A:447:ARG:HH11	1:A:668:ARG:HH12	1.47	0.62
2:C:211:PHE:HE2	2:C:228:ARG:HB2	1.65	0.62
3:D:1009:TYR:HD2	3:D:1090:MET:HE3	1.65	0.61
1:B:1193:TRP:HD1	1:B:1194:PHE:H	1.47	0.61
1:B:1060:ASP:OD2	1:B:1062:ARG:NH1	2.34	0.61
1:A:377:CYS:SG	1:A:378:THR:N	2.74	0.60
2:C:489:ASP:N	2:C:489:ASP:OD1	2.33	0.60
1:A:1203:GLY:HA2	1:A:1209:MET:HE3	1.84	0.60
3:D:1303:CYS:HB3	3:D:1307:CYS:SG	2.42	0.59
1:A:790:ASP:HB3	1:A:793:LEU:HD23	1.83	0.59
2:C:1036:ARG:NH1	2:C:1072:THR:O	2.35	0.59
2:C:427:VAL:HG12	2:C:438:THR:HG22	1.85	0.59
3:D:1286:HIS:CE1	3:D:1307:CYS:HB3	2.37	0.59
1:A:705:LYS:HD3	1:A:709:SER:HB2	1.85	0.58
1:A:16:ASP:OD2	1:A:29:HIS:NE2	2.36	0.58
1:A:337:ARG:NH1	1:A:349:ASP:OD2	2.37	0.58
1:A:588:ASN:OD1	1:A:591:THR:OG1	2.20	0.58
2:C:651:THR:OG1	2:C:652:LYS:N	2.37	0.57
1:B:792:GLN:NE2	1:B:1015:ASN:OD1	2.38	0.57
1:B:499:THR:HG1	1:B:519:SER:HG	1.44	0.56
2:C:523:THR:OG1	2:C:545:ASP:OD1	2.22	0.56
3:D:1047:ASP:OD1	3:D:1048:LEU:N	2.38	0.56
2:C:1107:CYS:O	2:C:1108:HIS:ND1	2.39	0.55
1:A:960:VAL:HG21	1:A:981:VAL:HG21	1.88	0.55
2:C:806:LYS:NZ	2:C:810:GLU:OE2	2.35	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:924:VAL:HG12	3:D:935:VAL:HG12	1.89	0.54
1:A:915:ALA:HB2	1:A:924:LEU:HB2	1.89	0.54
3:D:734:SER:OG	3:D:735:TYR:N	2.40	0.54
1:B:471:TRP:CZ3	1:B:490:LYS:HB3	2.44	0.53
1:B:1199:ASN:N	1:B:1199:ASN:OD1	2.41	0.53
3:D:845:ASN:O	3:D:845:ASN:ND2	2.35	0.53
1:B:243:PHE:HA	1:B:248:VAL:HA	1.90	0.53
1:B:144:ILE:O	1:B:178:LEU:N	2.41	0.53
1:B:669:VAL:HB	1:B:678:LEU:HD11	1.91	0.53
1:B:693:ASN:ND2	1:B:697:ASN:O	2.42	0.53
1:A:342:THR:OG1	1:A:346:THR:OG1	2.17	0.53
1:B:827:LYS:NZ	1:B:1004:GLU:OE1	2.37	0.53
1:B:759:ASP:OD1	1:B:760:ILE:N	2.42	0.53
2:C:83:HIS:CE1	2:C:85:MET:HB3	2.44	0.52
1:A:354:GLU:N	1:A:354:GLU:OE1	2.43	0.52
1:A:1142:VAL:HG23	1:A:1143:THR:HG23	1.90	0.52
1:B:358:LEU:HD23	1:B:366:LEU:HD11	1.91	0.52
1:B:1144:SER:HB2	1:B:1149:ARG:HA	1.91	0.52
3:D:543:ARG:NH1	3:D:676:GLN:OE1	2.43	0.52
1:A:1201:ARG:NH2	2:C:1099:SER:O	2.43	0.52
1:A:144:ILE:HD12	1:A:179:TYR:HB2	1.91	0.51
1:B:67:ALA:HB3	1:B:79:ALA:HB3	1.92	0.51
2:C:588:SER:OG	2:C:610:GLU:OE2	2.27	0.51
1:B:582:VAL:HG23	1:B:600:LEU:HB2	1.93	0.51
1:A:226:ALA:HA	1:A:244:TYR:HB2	1.92	0.51
1:A:1157:LEU:HD12	1:A:1208:HIS:HB3	1.92	0.51
2:C:691:TRP:NE1	2:C:707:ASP:OD1	2.34	0.51
1:A:273:VAL:HG22	1:A:285:THR:HG22	1.93	0.50
2:C:77:CYS:HG	2:C:92:THR:HG1	1.57	0.50
1:A:508:PRO:HD2	1:A:511:MET:HE2	1.91	0.50
1:A:1102:GLN:NE2	2:C:1113:MET:O	2.44	0.50
3:D:807:ILE:HG13	3:D:843:CYS:HB2	1.94	0.50
3:D:1034:ASP:N	3:D:1034:ASP:OD1	2.44	0.50
3:D:1303:CYS:SG	3:D:1305:THR:OG1	2.67	0.50
2:C:440:SER:OG	2:C:441:SER:N	2.45	0.49
1:A:995:SER:HB3	1:A:998:GLN:HG2	1.95	0.49
1:A:503:ASP:OD1	1:A:504:VAL:N	2.45	0.49
1:B:981:VAL:HG13	1:B:986:ARG:HD2	1.93	0.49
1:A:102:ASP:OD1	1:A:102:ASP:N	2.27	0.49
1:A:1115:SER:OG	1:A:1116:TYR:N	2.46	0.49
1:B:355:ARG:HH21	1:B:667:TRP:HE1	1.59	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:492:THR:HG21	1:B:538:TRP:CD1	2.47	0.49
1:B:1186:LYS:HA	1:B:1189:GLN:HG2	1.95	0.49
3:D:1246:SER:O	3:D:1248:CYS:N	2.46	0.49
1:A:16:ASP:OD1	1:A:16:ASP:N	2.38	0.48
3:D:839:ALA:HB2	3:D:876:TRP:HE1	1.78	0.48
1:A:620:PHE:HB3	1:A:636:LYS:HG2	1.95	0.48
1:A:266:GLN:HG3	1:A:270:LEU:HD21	1.96	0.48
3:D:736:ASP:O	3:D:737:ARG:HG2	2.14	0.48
1:B:472:GLU:OE1	1:B:474:ILE:N	2.47	0.47
2:C:35:CYS:SG	2:C:36:ARG:N	2.88	0.47
2:C:486:CYS:HB2	2:C:525:VAL:HG23	1.94	0.47
1:A:1164:THR:OG1	2:C:1129:GLN:OE1	2.31	0.47
1:B:462:CYS:HB2	1:B:501:VAL:HG23	1.97	0.47
1:B:627:THR:HG23	1:B:628:LYS:H	1.79	0.47
2:C:577:ILE:HG22	2:C:593:VAL:HG12	1.97	0.47
1:A:950:ARG:NH2	1:A:963:ALA:HB2	2.29	0.47
1:B:411:ARG:HD3	1:B:450:TRP:HH2	1.78	0.47
1:B:690:TRP:CZ3	1:B:700:CYS:HB2	2.50	0.47
1:A:378:THR:OG1	1:A:379:GLU:N	2.48	0.47
1:B:374:LEU:HD13	1:B:412:MET:HE1	1.96	0.47
1:B:688:ARG:CZ	1:B:703:ILE:HD12	2.45	0.47
3:D:1296:TRP:CH2	3:D:1300:GLN:HG3	2.50	0.47
1:B:1212:TRP:CD1	1:B:1220:PRO:HG3	2.50	0.46
1:A:502:THR:OG1	1:A:518:CYS:SG	2.72	0.46
3:D:1248:CYS:HB3	3:D:1260:CYS:HB3	1.63	0.46
1:B:404:SER:OG	1:B:448:VAL:O	2.23	0.46
2:C:27:ALA:HB3	2:C:48:ILE:HB	1.97	0.46
1:B:403:VAL:HG13	1:B:412:MET:HE3	1.97	0.46
1:A:1004:GLU:O	1:A:1008:ASN:ND2	2.45	0.45
1:B:521:ASP:N	1:B:521:ASP:OD1	2.49	0.45
1:B:330:TRP:HA	1:B:338:MET:HA	1.97	0.45
1:B:398:ASP:N	1:B:398:ASP:OD1	2.49	0.45
1:B:629:ASP:HB2	1:B:659:PHE:O	2.16	0.45
2:C:347:LEU:HD12	2:C:697:ILE:HD13	1.98	0.45
3:D:845:ASN:OD1	3:D:869:ASP:N	2.49	0.45
1:A:313:ILE:HD13	1:A:314:ILE:H	1.82	0.45
2:C:159:PHE:CE1	2:C:196:PRO:HB2	2.52	0.45
1:A:635:LEU:HD22	1:A:653:ILE:HG13	1.99	0.45
1:B:1226:CYS:SG	1:B:1227:LYS:N	2.89	0.45
3:D:677:MET:HE2	3:D:681:GLY:HA2	1.98	0.45
2:C:249:GLN:N	2:C:249:GLN:OE1	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:443:GLN:HG2	2:C:467:GLY:O	2.16	0.45
3:D:1274:VAL:HG22	3:D:1289:HIS:NE2	2.32	0.45
3:D:805:HIS:CD2	3:D:842:GLY:HA3	2.52	0.45
1:A:467:THR:HG22	1:A:495:VAL:HG12	1.98	0.45
1:B:452:HIS:ND1	1:B:453:PRO:HD2	2.31	0.45
1:B:472:GLU:OE1	1:B:473:GLU:N	2.50	0.45
1:B:806:TYR:CD1	1:B:822:VAL:HG11	2.52	0.45
1:B:434:CYS:SG	1:B:435:THR:N	2.90	0.45
3:D:1192:ARG:HH21	3:D:1196:PRO:HG3	1.81	0.45
2:C:1099:SER:OG	2:C:1100:ARG:N	2.50	0.44
1:A:234:PRO:O	1:A:237:HIS:NE2	2.50	0.44
3:D:636:SER:OG	3:D:637:GLY:N	2.50	0.44
1:A:53:LEU:HD21	1:A:56:ILE:HG12	1.99	0.44
1:B:145:TRP:HA	1:B:178:LEU:H	1.82	0.44
1:A:1032:MET:HG3	1:A:1049:MET:HE2	1.99	0.44
2:C:463:LYS:HE3	2:C:463:LYS:HB3	1.87	0.44
2:C:426:ASP:OD1	2:C:427:VAL:N	2.51	0.44
1:B:140:PHE:HA	1:B:183:GLN:HA	2.00	0.44
1:B:693:ASN:ND2	1:B:697:ASN:OD1	2.35	0.44
3:D:709:LEU:HD13	3:D:742:TRP:CE3	2.53	0.44
1:B:455:PHE:CD2	1:B:513:LEU:HD12	2.53	0.44
1:B:1029:VAL:HG21	1:B:1062:ARG:HH12	1.83	0.43
2:C:41:VAL:HG13	2:C:52:TYR:HB2	2.01	0.43
1:A:950:ARG:NH2	1:A:959:CYS:O	2.45	0.43
1:B:781:ARG:O	1:B:784:ILE:HG12	2.19	0.43
2:C:275:ASP:OD1	2:C:276:HIS:N	2.51	0.43
1:A:297:ASP:HB3	1:A:314:ILE:HG13	2.00	0.43
3:D:556:PRO:HG3	3:D:645:GLY:HA2	2.00	0.43
3:D:678:ASP:OD1	3:D:680:TYR:N	2.51	0.43
1:A:665:GLN:HE21	1:A:665:GLN:HB3	1.63	0.43
1:A:973:ASP:OD1	1:A:973:ASP:N	2.46	0.43
1:B:1155:CYS:SG	1:B:1156:ALA:N	2.91	0.43
3:D:1065:HIS:CD2	3:D:1066:PRO:HD2	2.53	0.43
1:B:242:SER:O	1:B:249:ALA:N	2.41	0.43
1:B:629:ASP:OD1	1:B:629:ASP:N	2.51	0.43
3:D:767:TRP:CZ3	3:D:776:LEU:HB2	2.53	0.43
1:A:819:GLY:HA2	1:A:822:VAL:HG12	2.00	0.43
1:A:1016:LEU:HD23	1:A:1016:LEU:HA	1.85	0.43
1:B:1107:CYS:HB2	3:D:1239:ILE:HD13	2.01	0.43
2:C:53:ALA:HB3	2:C:60:VAL:HG23	2.00	0.43
3:D:927:SER:OG	3:D:928:ILE:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:LEU:HD12	1:A:392:ILE:HD11	1.99	0.42
1:A:442:SER:HB3	1:A:465:ASP:OD2	2.18	0.42
2:C:266:HIS:CE1	2:C:282:ASP:HB3	2.55	0.42
1:B:691:LYS:HE3	1:B:691:LYS:HB2	1.84	0.42
1:B:1098:LYS:HE3	3:D:1306:GLY:O	2.20	0.42
3:D:1257:CYS:SG	3:D:1258:ALA:N	2.92	0.42
1:B:188:LEU:N	1:B:203:GLY:O	2.51	0.42
1:B:620:PHE:HB3	1:B:636:LYS:HG2	2.02	0.42
2:C:318:SER:OG	2:C:319:LYS:N	2.53	0.42
3:D:690:ASP:O	3:D:691:ARG:HG2	2.19	0.42
1:A:923:ASN:O	1:A:927:VAL:HG23	2.20	0.42
2:C:221:GLY:O	2:C:237:MET:N	2.44	0.42
3:D:863:LYS:HE2	3:D:863:LYS:HB3	1.93	0.42
1:B:695:MET:O	1:B:697:ASN:ND2	2.53	0.42
1:B:1073:LEU:HD23	1:B:1073:LEU:HA	1.89	0.42
2:C:708:ASP:OD1	2:C:708:ASP:N	2.51	0.42
2:C:987:LEU:HD13	2:C:987:LEU:HA	1.88	0.42
1:B:518:CYS:HB2	1:B:524:VAL:HG22	2.01	0.41
3:D:555:ARG:NH1	3:D:635:ASP:HB3	2.36	0.41
3:D:1065:HIS:CG	3:D:1066:PRO:HD2	2.55	0.41
1:B:941:LEU:HD12	1:B:942:TRP:N	2.34	0.41
2:C:143:GLY:HA2	2:C:166:VAL:HG23	2.02	0.41
2:C:255:ALA:HB2	2:C:272:MET:HB3	2.01	0.41
1:A:857:ASN:O	1:A:861:ILE:HG12	2.21	0.41
1:A:122:ASN:HB3	1:A:125:ASP:O	2.21	0.41
2:C:466:SER:N	2:C:489:ASP:OD2	2.48	0.41
1:A:69:TYR:CD2	1:A:77:LEU:HD23	2.57	0.40
1:B:693:ASN:HD21	1:B:697:ASN:CG	2.27	0.40
2:C:414:ARG:H	2:C:414:ARG:HG3	1.75	0.40
2:C:465:HIS:CG	2:C:487:SER:HG	2.39	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1089/1113 (98%)	1034 (95%)	55 (5%)	0	100	100
1	B	1070/1113 (96%)	1031 (96%)	39 (4%)	0	100	100
2	C	893/905 (99%)	864 (97%)	29 (3%)	0	100	100
3	D	621/633 (98%)	584 (94%)	36 (6%)	1 (0%)	43	74
All	All	3673/3764 (98%)	3513 (96%)	159 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	D	1247	HIS

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	835/977 (86%)	805 (96%)	30 (4%)	31	57
1	B	580/977 (59%)	544 (94%)	36 (6%)	16	44
2	C	680/785 (87%)	642 (94%)	38 (6%)	19	46
3	D	462/537 (86%)	432 (94%)	30 (6%)	15	42
All	All	2557/3276 (78%)	2423 (95%)	134 (5%)	22	48

All (134) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	20	VAL
1	A	70	LEU
1	A	73	ASP
1	A	92	LEU
1	A	102	ASP
1	A	109	VAL

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Mol	Chain	Res	Type
1	A	111	LYS
1	A	145	TRP
1	A	215	ASN
1	A	287	THR
1	A	313	ILE
1	A	314	ILE
1	A	347	MET
1	A	353	PHE
1	A	403	VAL
1	A	414	THR
1	A	466	ARG
1	A	517	THR
1	A	535	LEU
1	A	544	ILE
1	A	616	LEU
1	A	634	THR
1	A	665	GLN
1	A	778	GLN
1	A	911	LEU
1	A	1019	ILE
1	A	1100	LEU
1	A	1187	LEU
1	A	1197	CYS
1	B	353	PHE
1	B	366	LEU
1	B	398	ASP
1	B	405	PHE
1	B	414	THR
1	B	449	THR
1	B	455	PHE
1	B	502	THR
1	B	513	LEU
1	B	540	LEU
1	B	581	LYS
1	B	616	LEU
1	B	627	THR
1	B	655	ILE
1	B	662	HIS
1	B	683	ASP
1	B	697	ASN
1	B	703	ILE
1	B	769	LEU

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Mol	Chain	Res	Type
1	B	774	LEU
1	B	807	THR
1	B	878	VAL
1	B	899	LEU
1	B	905	ARG
1	B	929	MET
1	B	941	LEU
1	B	999	LEU
1	B	1058	LEU
1	B	1060	ASP
1	B	1062	ARG
1	B	1074	LEU
1	B	1105	VAL
1	B	1144	SER
1	B	1197	CYS
1	B	1200	CYS
1	B	1210	LEU
2	C	21	MET
2	C	41	VAL
2	C	43	VAL
2	C	117	THR
2	C	182	THR
2	C	194	ARG
2	C	231	MET
2	C	237	MET
2	C	258	LYS
2	C	266	HIS
2	C	282	ASP
2	C	284	ARG
2	C	303	ILE
2	C	397	LEU
2	C	463	LYS
2	C	472	VAL
2	C	489	ASP
2	C	527	ASP
2	C	528	VAL
2	C	534	HIS
2	C	542	CYS
2	C	612	ASN
2	C	631	VAL
2	C	651	THR
2	C	652	LYS

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Mol	Chain	Res	Type
2	C	654	VAL
2	C	715	LYS
2	C	727	ILE
2	C	784	VAL
2	C	792	LEU
2	C	987	LEU
2	C	1031	LEU
2	C	1094	LYS
2	C	1121	VAL
2	C	1126	VAL
2	C	1138	GLN
2	C	1144	LEU
2	C	1159	GLU
3	D	539	THR
3	D	543	ARG
3	D	647	VAL
3	D	663	THR
3	D	691	ARG
3	D	697	ASP
3	D	712	HIS
3	D	762	VAL
3	D	770	HIS
3	D	774	LEU
3	D	807	ILE
3	D	845	ASN
3	D	863	LYS
3	D	873	ASP
3	D	896	VAL
3	D	913	LEU
3	D	922	TRP
3	D	942	VAL
3	D	943	THR
3	D	954	VAL
3	D	1034	ASP
3	D	1041	LEU
3	D	1088	LEU
3	D	1194	LEU
3	D	1251	GLU
3	D	1257	CYS
3	D	1267	CYS
3	D	1270	CYS
3	D	1284	CYS

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Mol	Chain	Res	Type
3	D	1312	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	HIS
1	A	419	GLN
1	A	441	HIS
1	A	534	ASN
1	A	542	HIS
1	A	778	GLN
1	A	857	ASN
1	A	1000	ASN
1	A	1043	GLN
1	A	1208	HIS
1	B	792	GLN
1	B	801	HIS
1	B	817	ASN
1	B	1043	GLN
2	C	83	HIS
2	C	124	ASN
2	C	185	ASN
2	C	307	HIS
2	C	686	HIS
3	D	753	HIS
3	D	1065	HIS
3	D	1078	HIS
3	D	1236	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates ⓘ

There are no 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	C	5
3	D	5
1	A	4
1	B	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	830:PRO	C	985:PHE	N	43.91
1	D	1099:SER	C	1178:ASP	N	27.16
1	C	730:GLY	C	779:GLY	N	22.41
1	A	710:PRO	C	757:GLU	N	21.53
1	A	1166:VAL	C	1183:LYS	N	19.23
1	A	1116:TYR	C	1141:LYS	N	16.40
1	C	360:VAL	C	389:PHE	N	16.12
1	D	957:SER	C	994:ASP	N	14.28
1	D	581:THR	C	629:SER	N	11.01
1	D	820:PRO	C	825:ASP	N	10.53
1	B	148:CYS	C	174:VAL	N	9.30
1	C	499:VAL	C	508:GLY	N	8.98
1	B	1116:TYR	C	1141:LYS	N	8.72
1	C	404:PRO	C	410:MET	N	4.68
1	A	379:GLU	C	386:MET	N	4.28
1	D	526:TYR	C	527:GLY	N	3.26

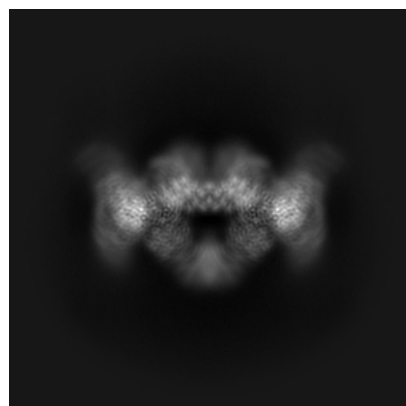
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-47276. These allow visual inspection of the internal detail of the map and identification of artifacts.

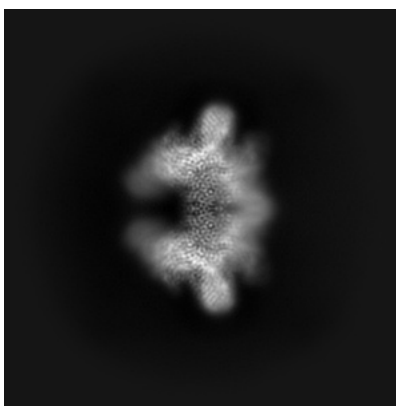
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

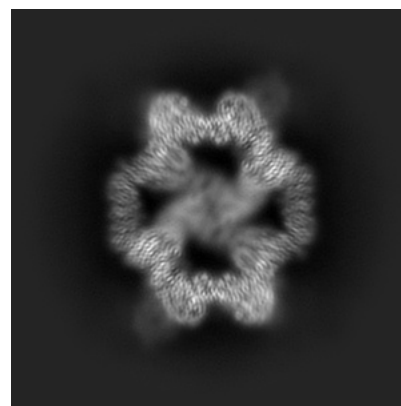
#### 6.1.1 Primary map



X

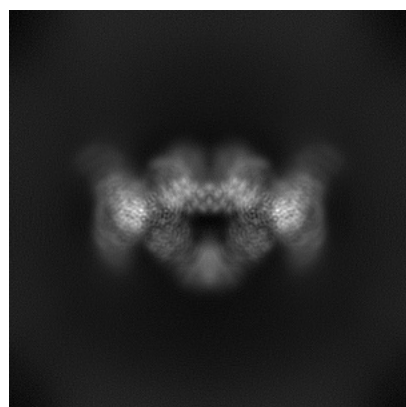


Y

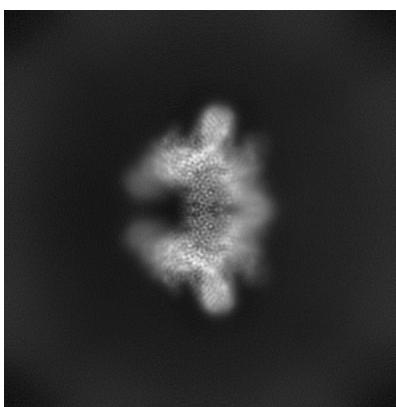


Z

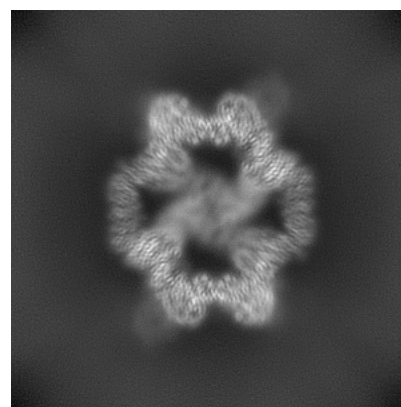
#### 6.1.2 Raw map



X



Y

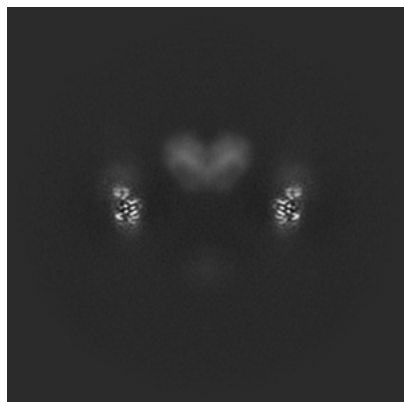


Z

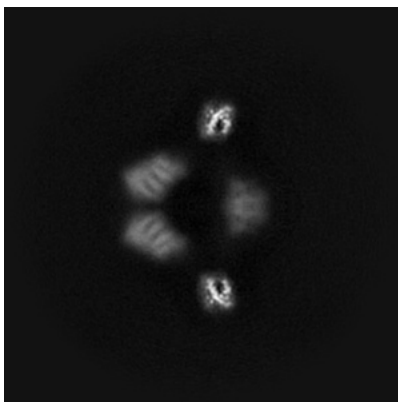
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

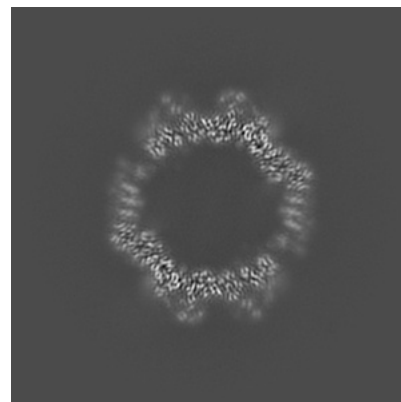
### 6.2.1 Primary map



X Index: 240

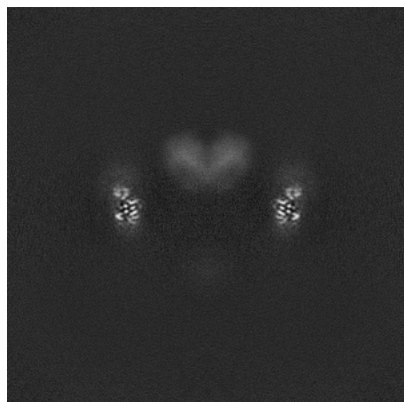


Y Index: 240

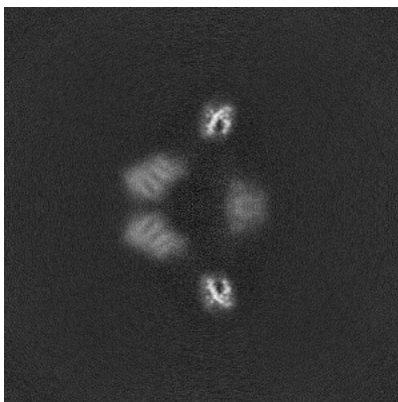


Z Index: 240

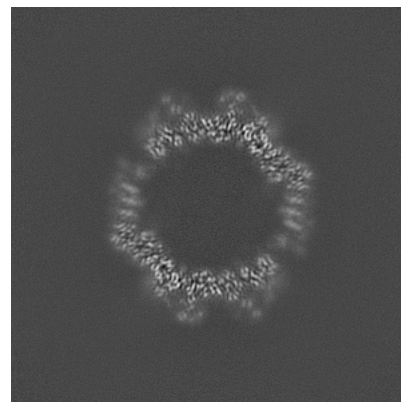
### 6.2.2 Raw map



X Index: 240



Y Index: 240

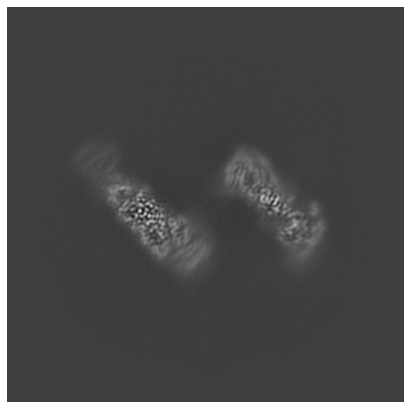


Z Index: 240

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

## 6.3 Largest variance slices [i](#)

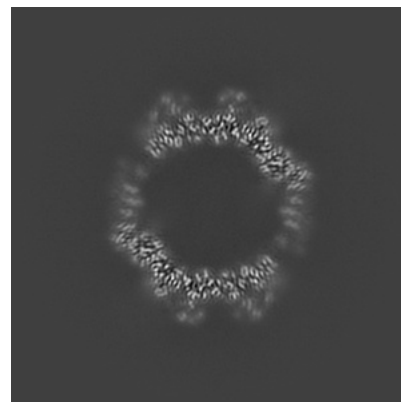
### 6.3.1 Primary map



X Index: 184

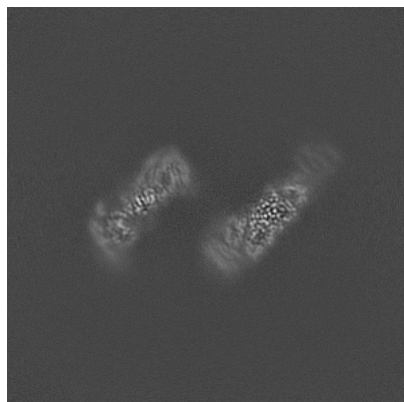


Y Index: 201

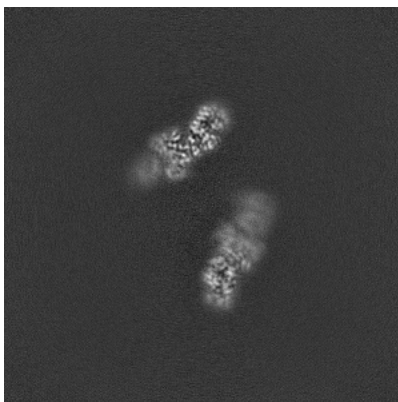


Z Index: 238

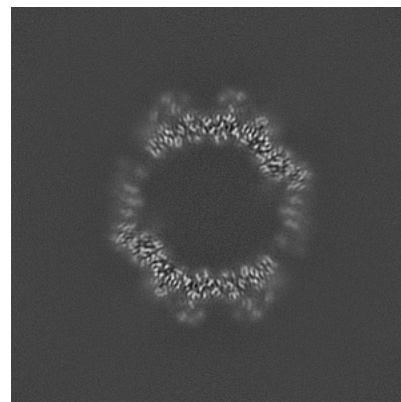
### 6.3.2 Raw map



X Index: 296



Y Index: 279

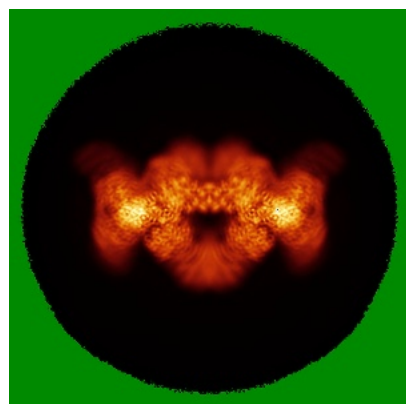


Z Index: 238

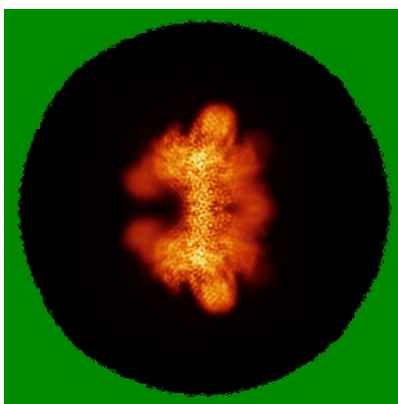
The images above show the largest variance slices of the map in three orthogonal directions.

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

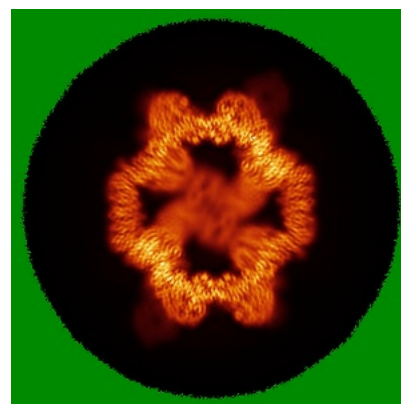
### 6.4.1 Primary map



X

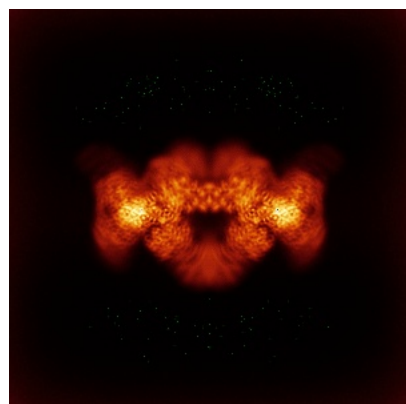


Y

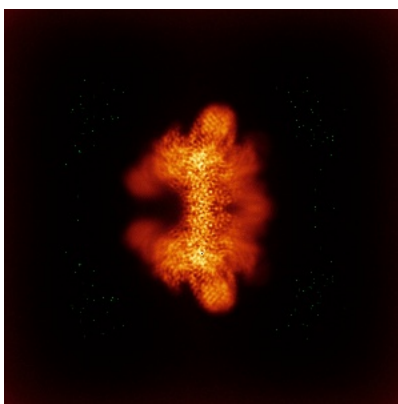


Z

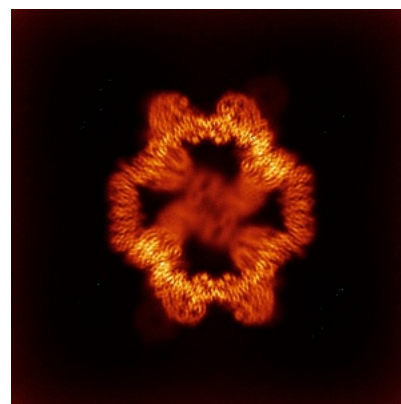
### 6.4.2 Raw map



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.



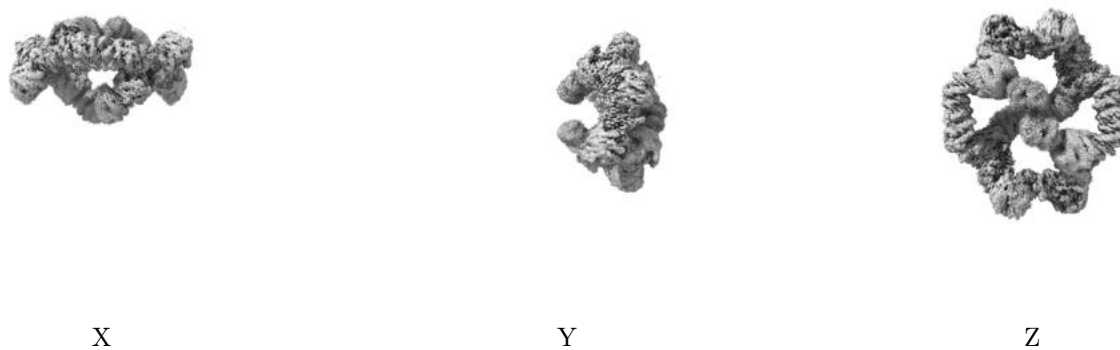
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.06. 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.



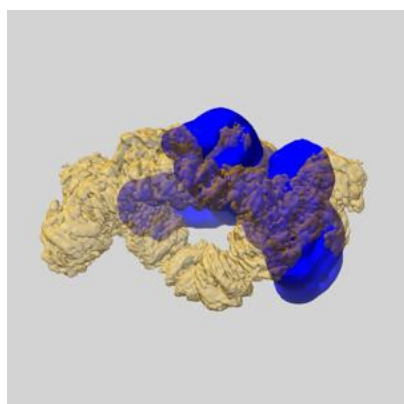
## 6.6 Mask visualisation [i](#)

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

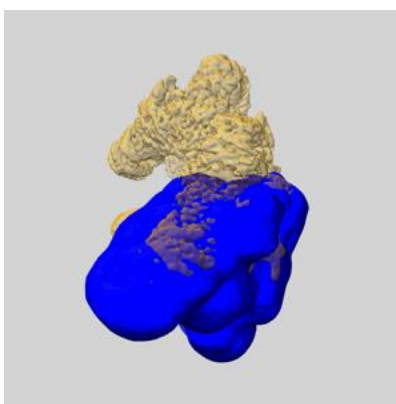
A mask typically either:

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

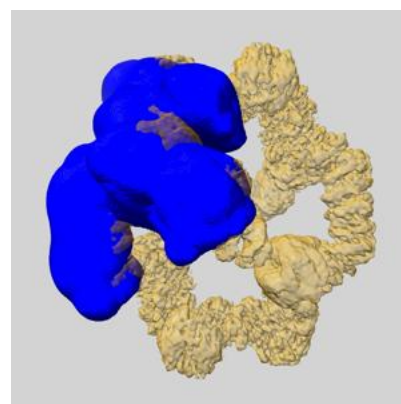
### 6.6.1 emd\_47276\_msk\_5.map [i](#)



X

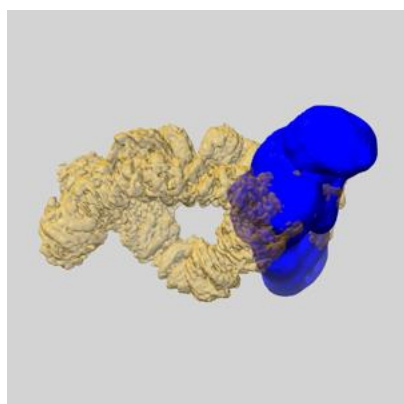


Y

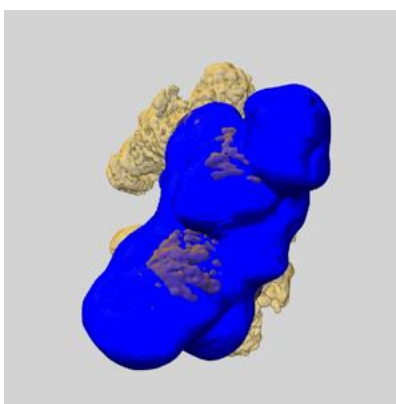


Z

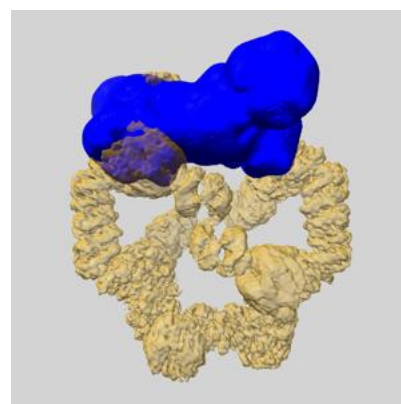
### 6.6.2 emd\_47276\_msk\_4.map [i](#)



X

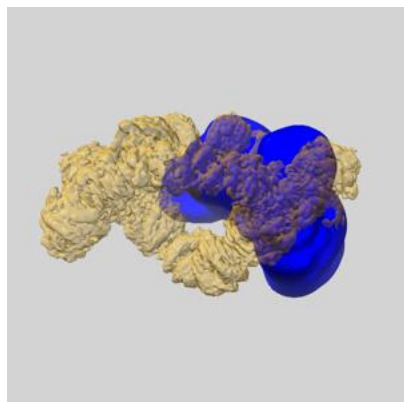


Y

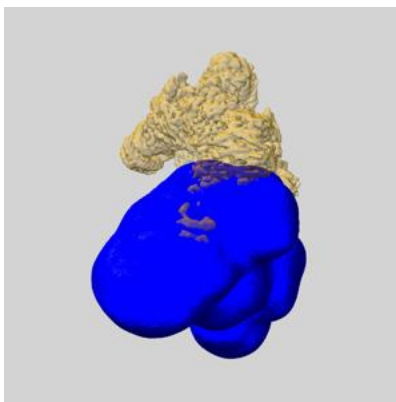


Z

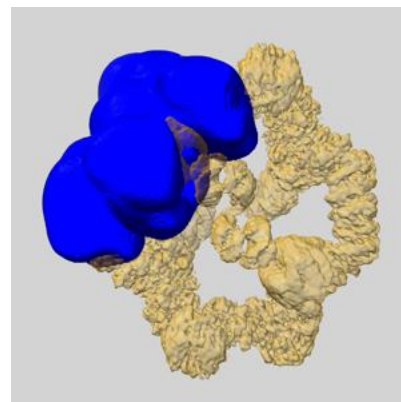
### 6.6.3 emd\_47276\_msk\_3.map [i](#)



X

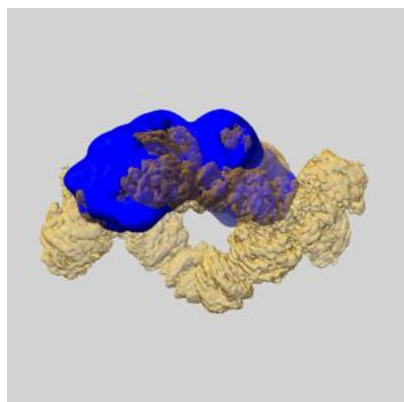


Y

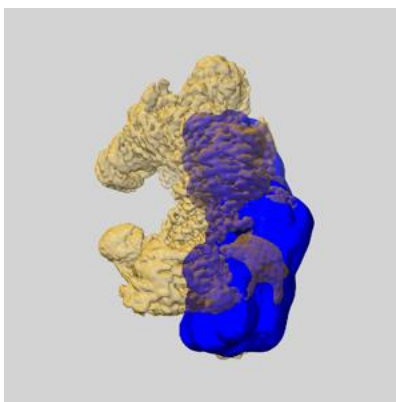


Z

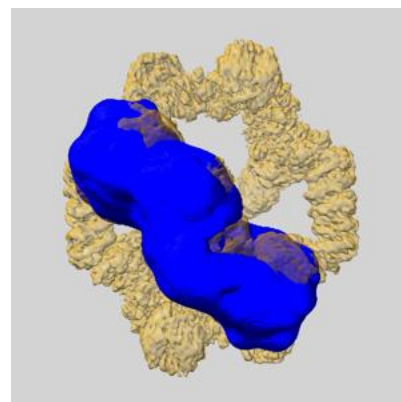
### 6.6.4 emd\_47276\_msk\_2.map [i](#)



X

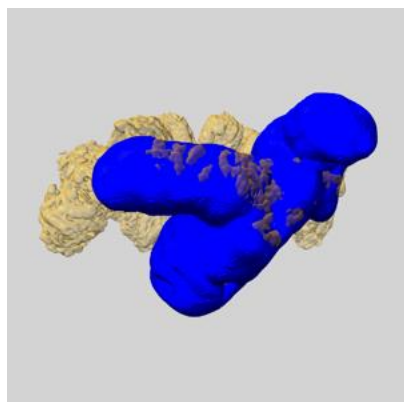


Y

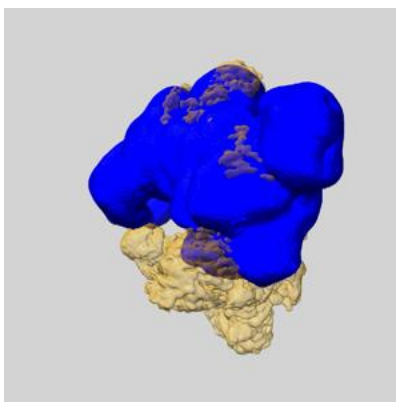


Z

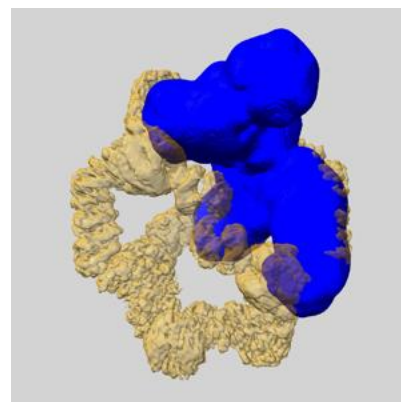
### 6.6.5 emd\_47276\_msk\_1.map [i](#)



X



Y

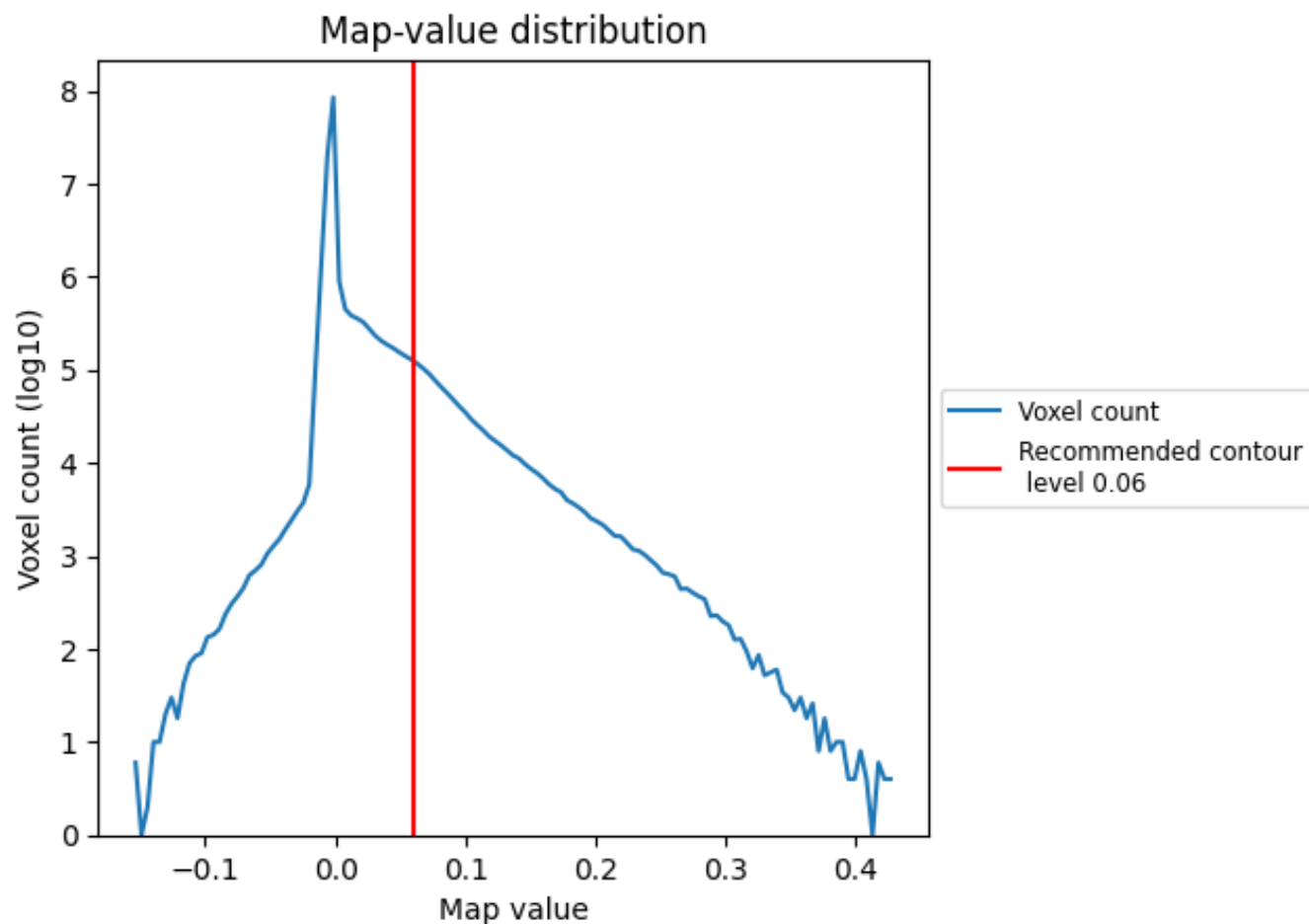


Z

## 7 Map analysis [i](#)

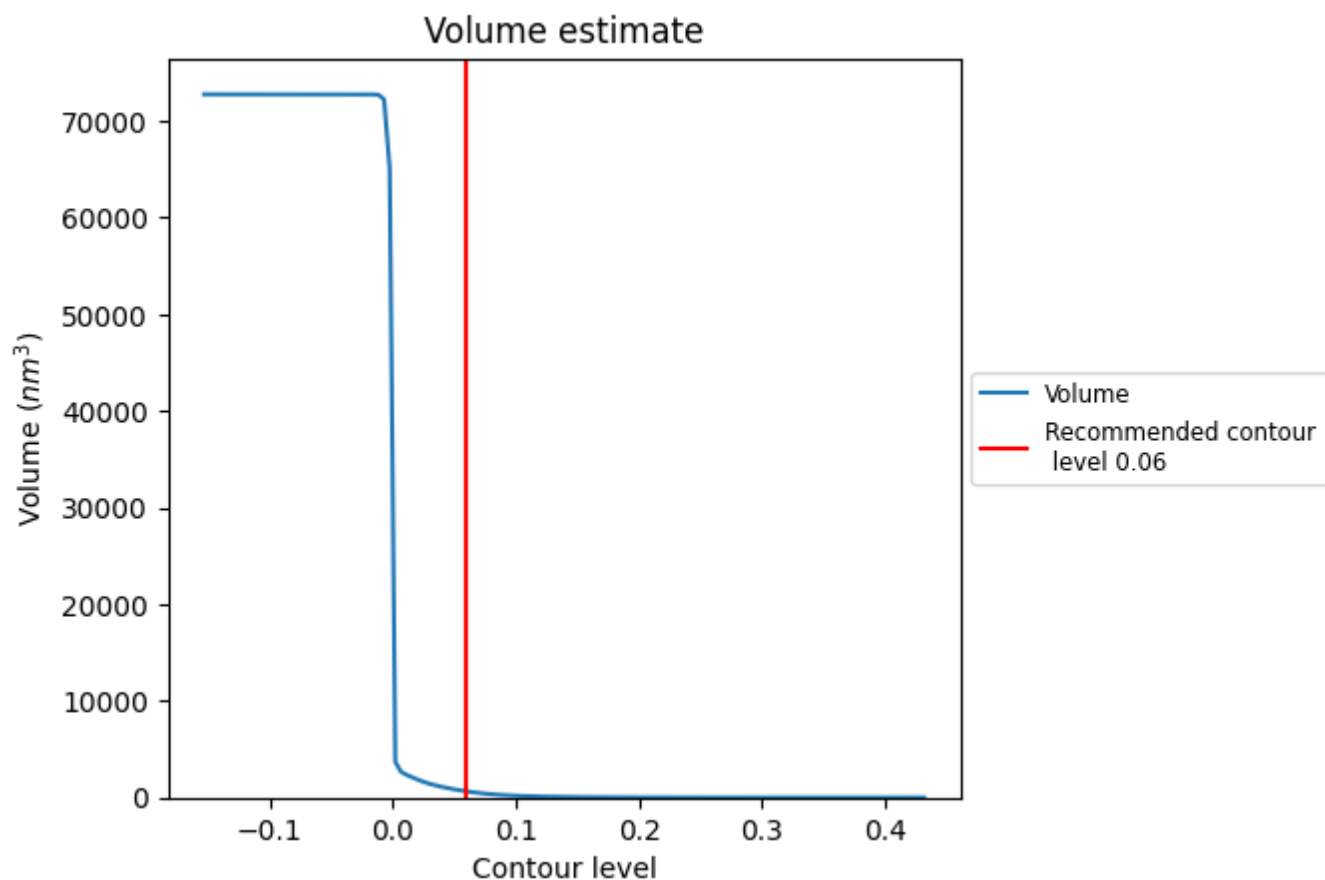
This section contains the results of statistical analysis of the map.

### 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.

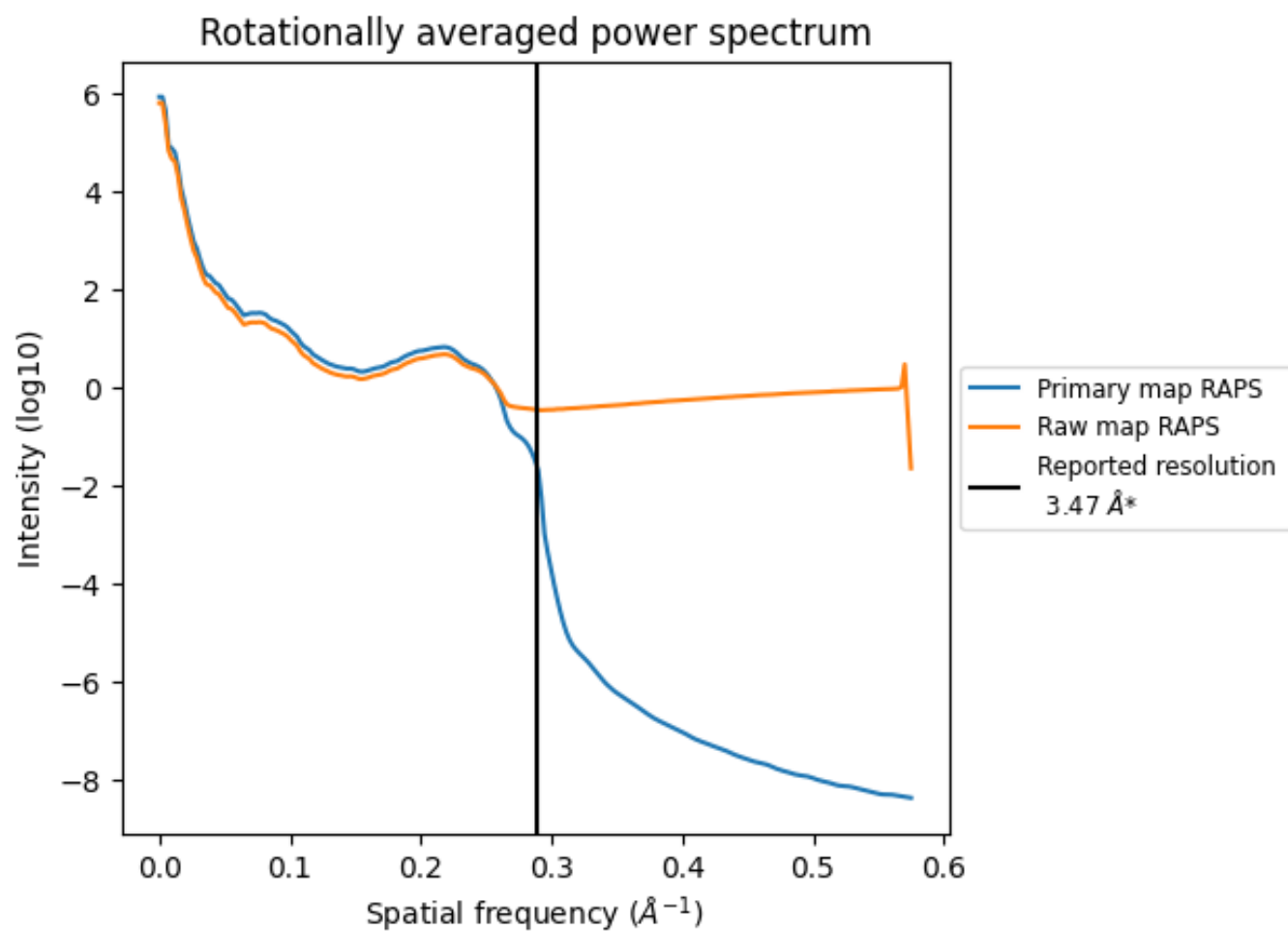
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 639  $\text{nm}^3$ ; this corresponds to an approximate mass of 577 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 ⓘ

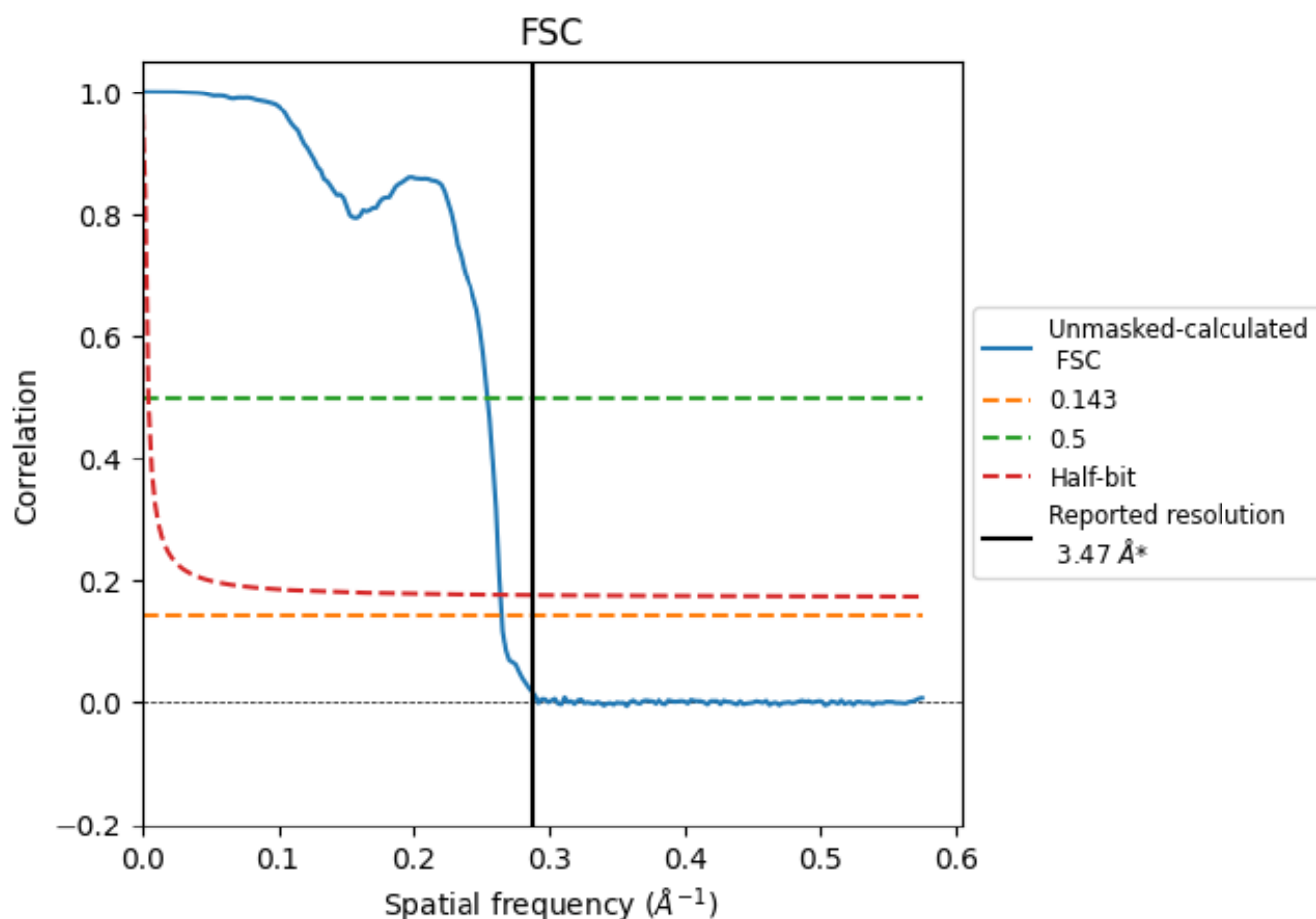


\*Reported resolution corresponds to spatial frequency of 0.288 Å<sup>-1</sup>

## 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 0.288  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

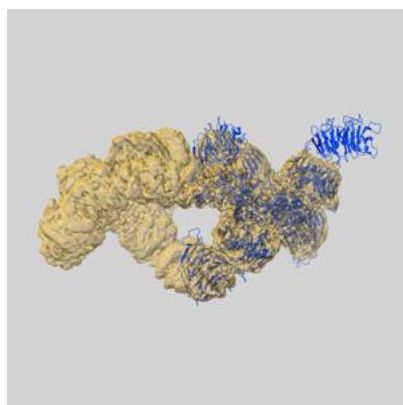
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.47	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.77	3.92	3.78

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

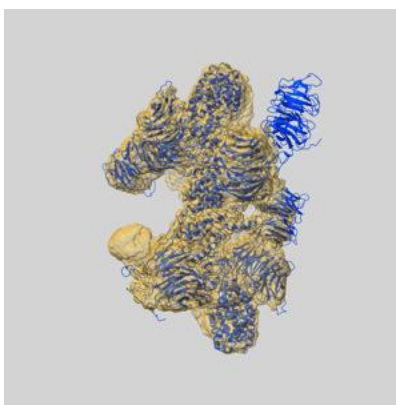
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-47276 and PDB model 9DX0. Per-residue inclusion information can be found in section [3](#) on page [4](#).

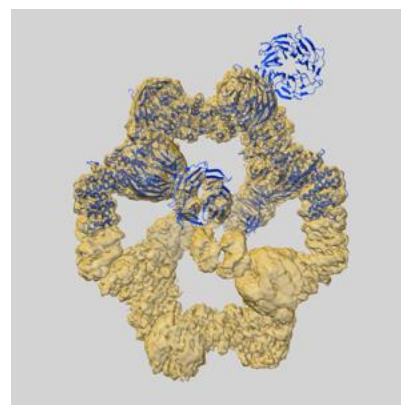
### 9.1 Map-model overlay [i](#)



X



Y

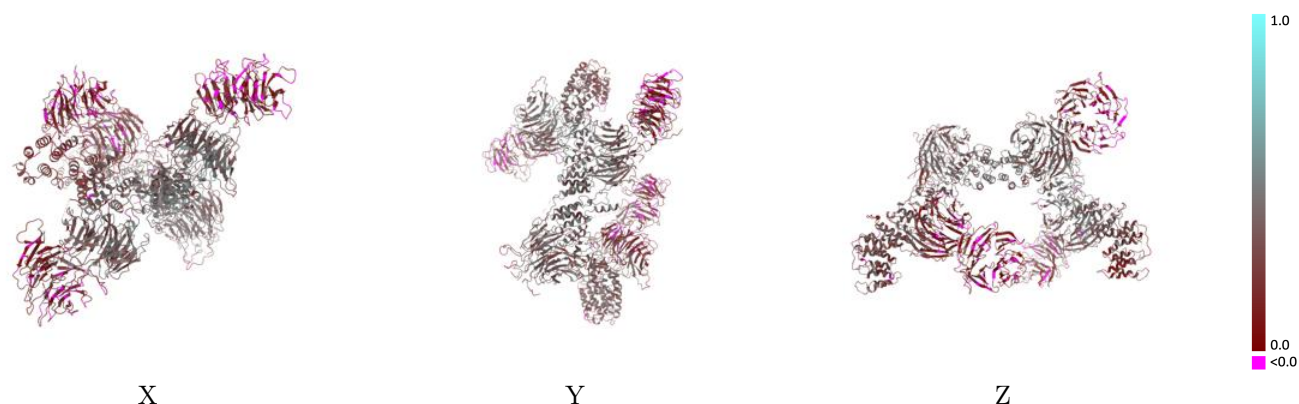


Z

The images above show the 3D surface view of the map at the recommended contour level 0.06 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.

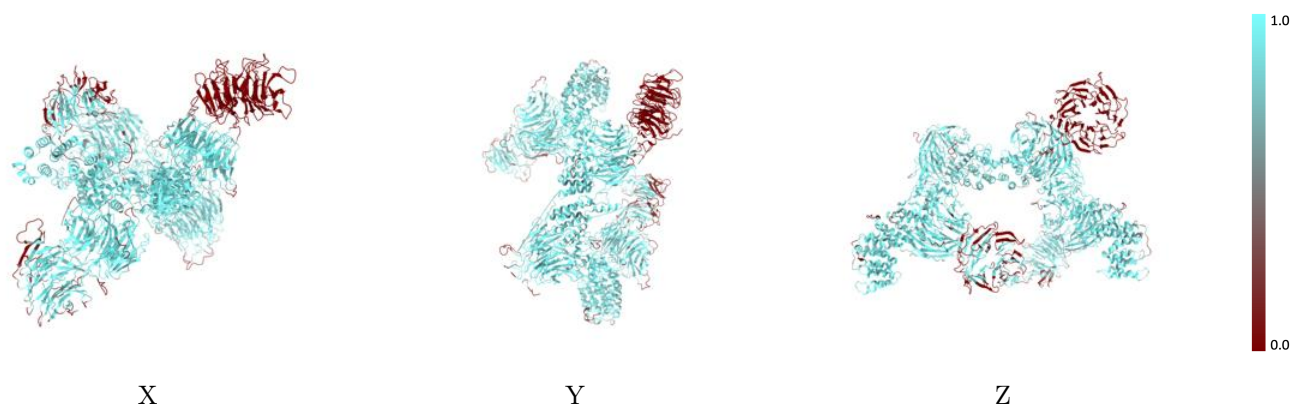


## 9.2 Q-score mapped to coordinate model [i](#)



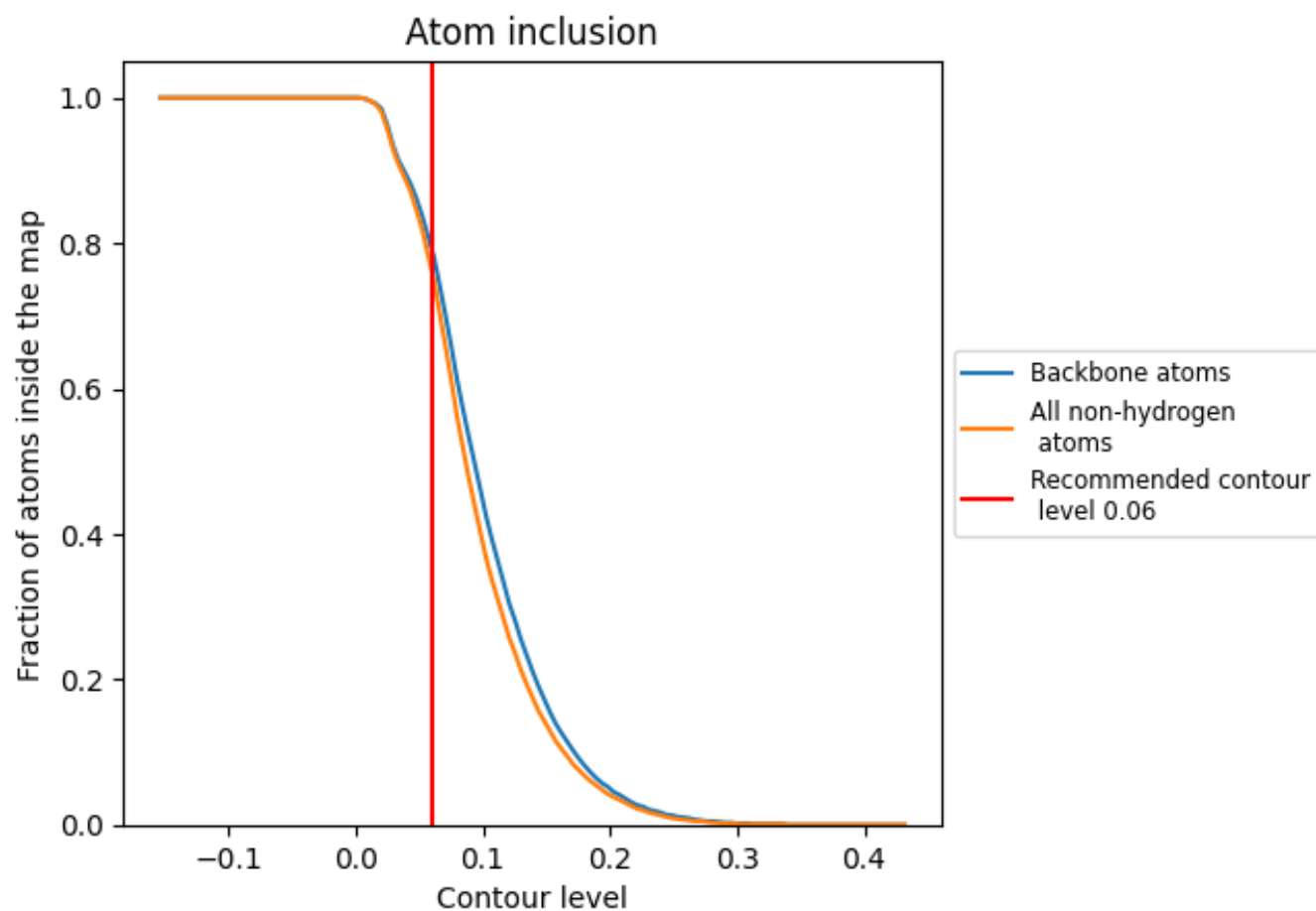
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.06).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 79% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.7610	<div></div> 0.3020
A	<div></div> 0.8350	<div></div> 0.3000
B	<div></div> 0.7840	<div></div> 0.2420
C	<div></div> 0.5790	<div></div> 0.3170
D	<div></div> 0.8550	<div></div> 0.3780

