



Full wwPDB EM Validation Report ⓘ

May 11, 2026 – 11:04 PM EDT

PDB ID : 9DX2 / pdb_00009dx2
EMDB ID : EMD-47278
Title : Human GATOR2 complex - CASTOR1 bound state
Authors : Wranik, M.; Rogala, K.B.
Deposited on : 2024-10-10
Resolution : 3.40 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

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:

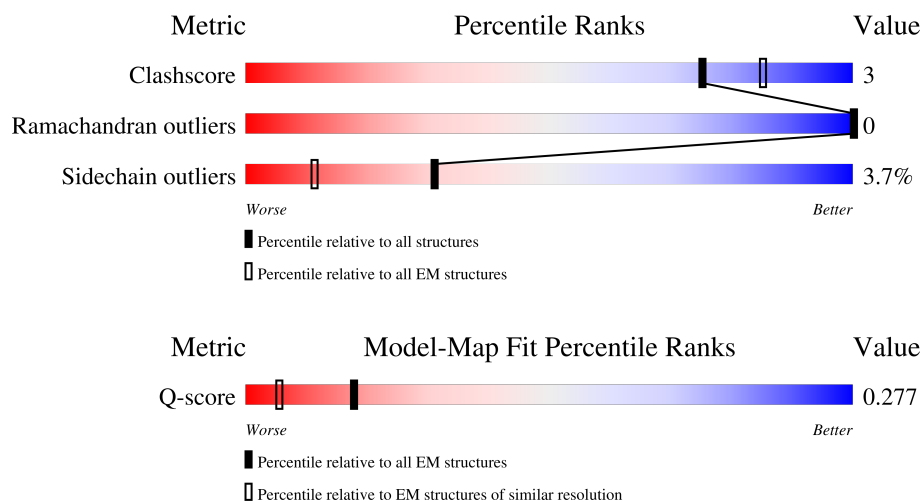
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.40 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	14717 (2.90 - 3.90)

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	A	786	 93% 7%
1	B	786	 87% 9% . .
2	C	570	 91% 9%
3	D	545	 92% 8%

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Mol	Chain	Length	Quality of chain
4	E	328	<div><div></div><div>11%</div><div>92%</div><div>7%</div><div></div></div>
4	F	328	<div><div></div><div>34%</div><div>83%</div><div>12%</div><div></div><div></div></div>
4	G	328	<div><div></div><div>67%</div><div>77%</div><div>9%</div><div>13%</div><div></div></div>
5	H	1042	<div><div></div><div>14%</div><div>88%</div><div>11%</div><div></div><div></div></div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 34252 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	786	Total	C	N	O	S	0	0
			5894	3721	1057	1070	46		
1	B	761	Total	C	N	O	S	0	0
			5582	3526	991	1021	44		

- Molecule 2 is a protein called Nucleoporin SEH1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	570	Total	C	N	O	S	0	0
			4367	2749	780	799	39		

- Molecule 3 is a protein called Protein SEC13 homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	545	Total	C	N	O	S	0	0
			4010	2539	714	735	22		

- Molecule 4 is a protein called Cytosolic arginine sensor for mTORC1 subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	328	Total	C	N	O	S	0	0
			2477	1593	413	464	7		
4	F	315	Total	C	N	O	S	0	0
			2279	1480	377	414	8		
4	G	284	Total	C	N	O	S	0	0
			2084	1356	347	375	6		

- Molecule 5 is a protein called GATOR2 complex protein WDR24.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	H	1042	Total	C	N	O	S	0	0
			7559	4822	1334	1351	52		

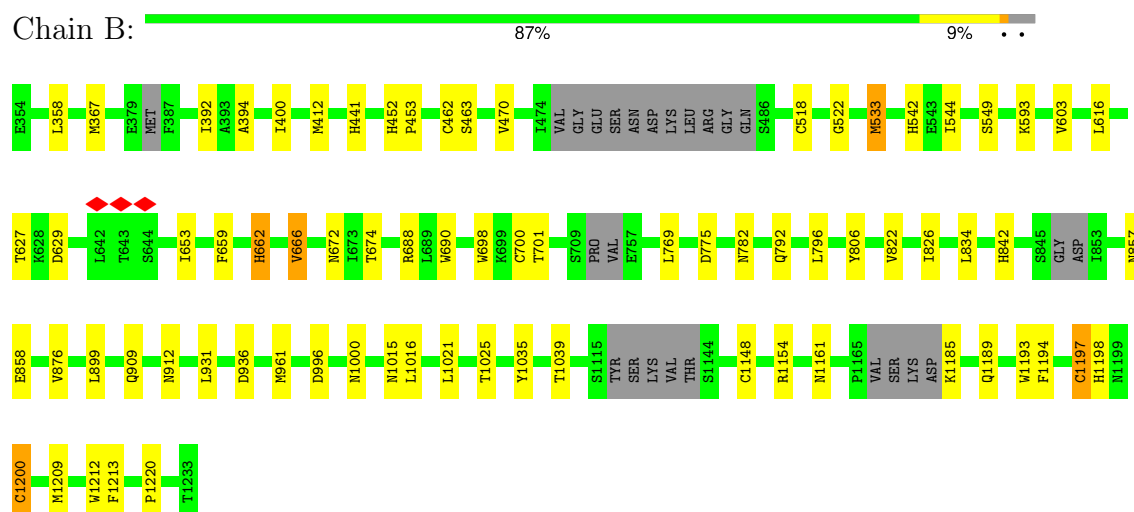
3 Residue-property plots

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.

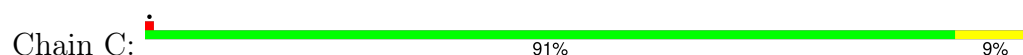
• Molecule 1: GATOR2 complex protein MIOS

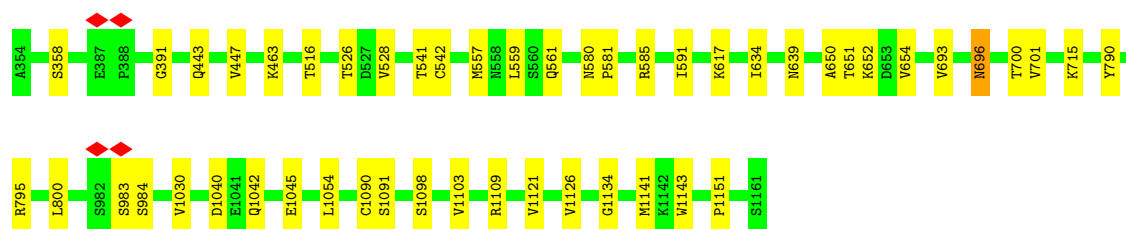


• Molecule 1: GATOR2 complex protein MIOS

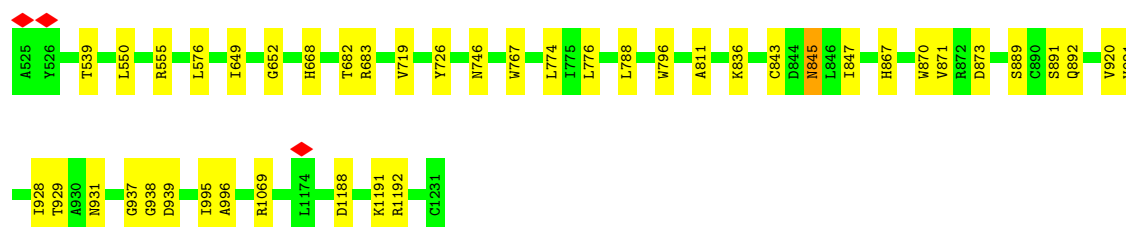


• Molecule 2: Nucleoporin SEH1





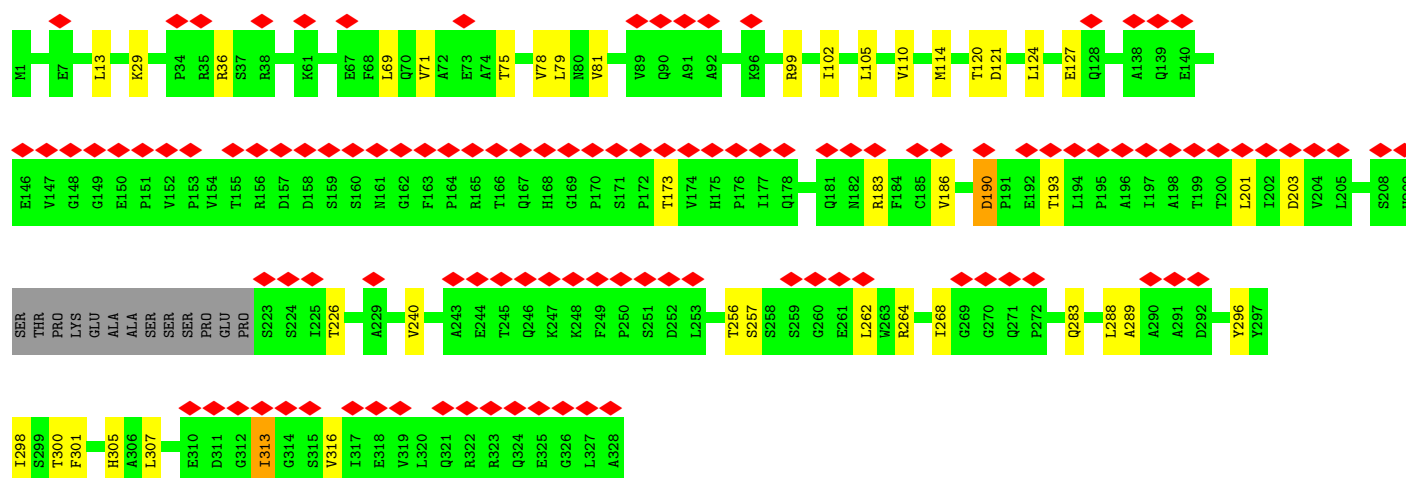
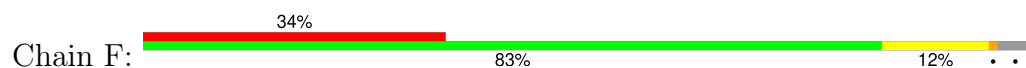
• Molecule 3: Protein SEC13 homolog



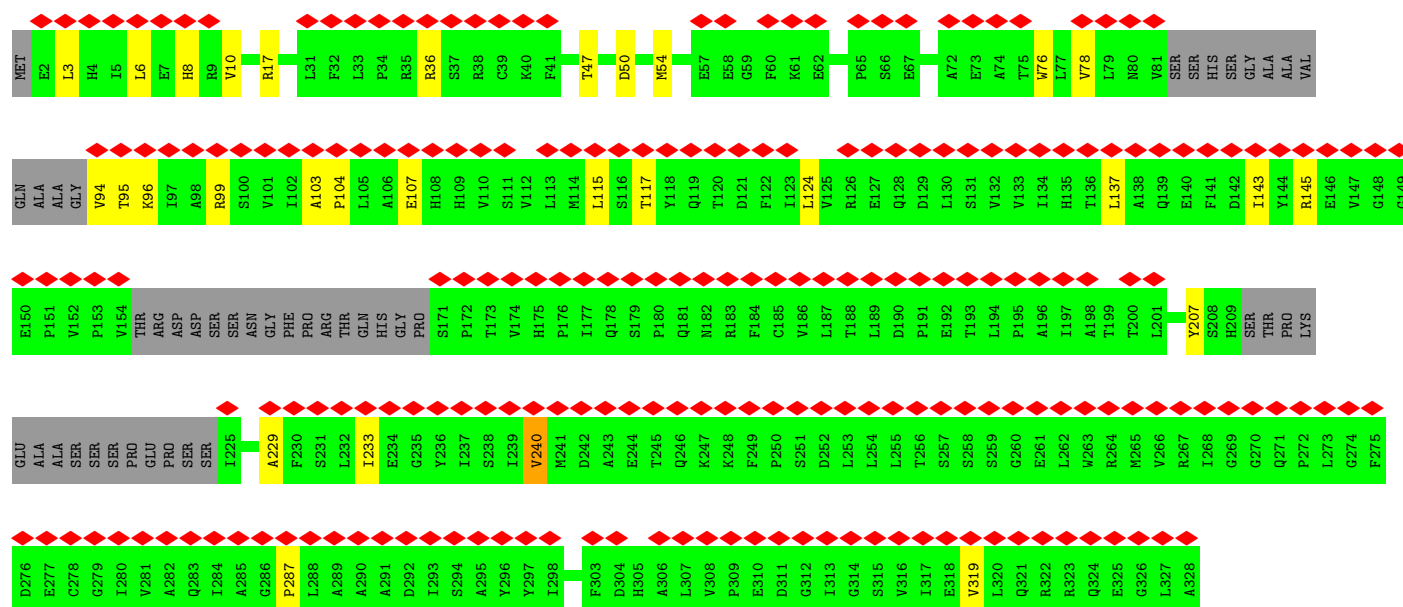
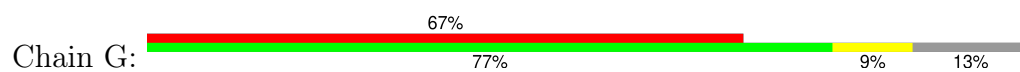
• Molecule 4: Cytosolic arginine sensor for mTORC1 subunit 1



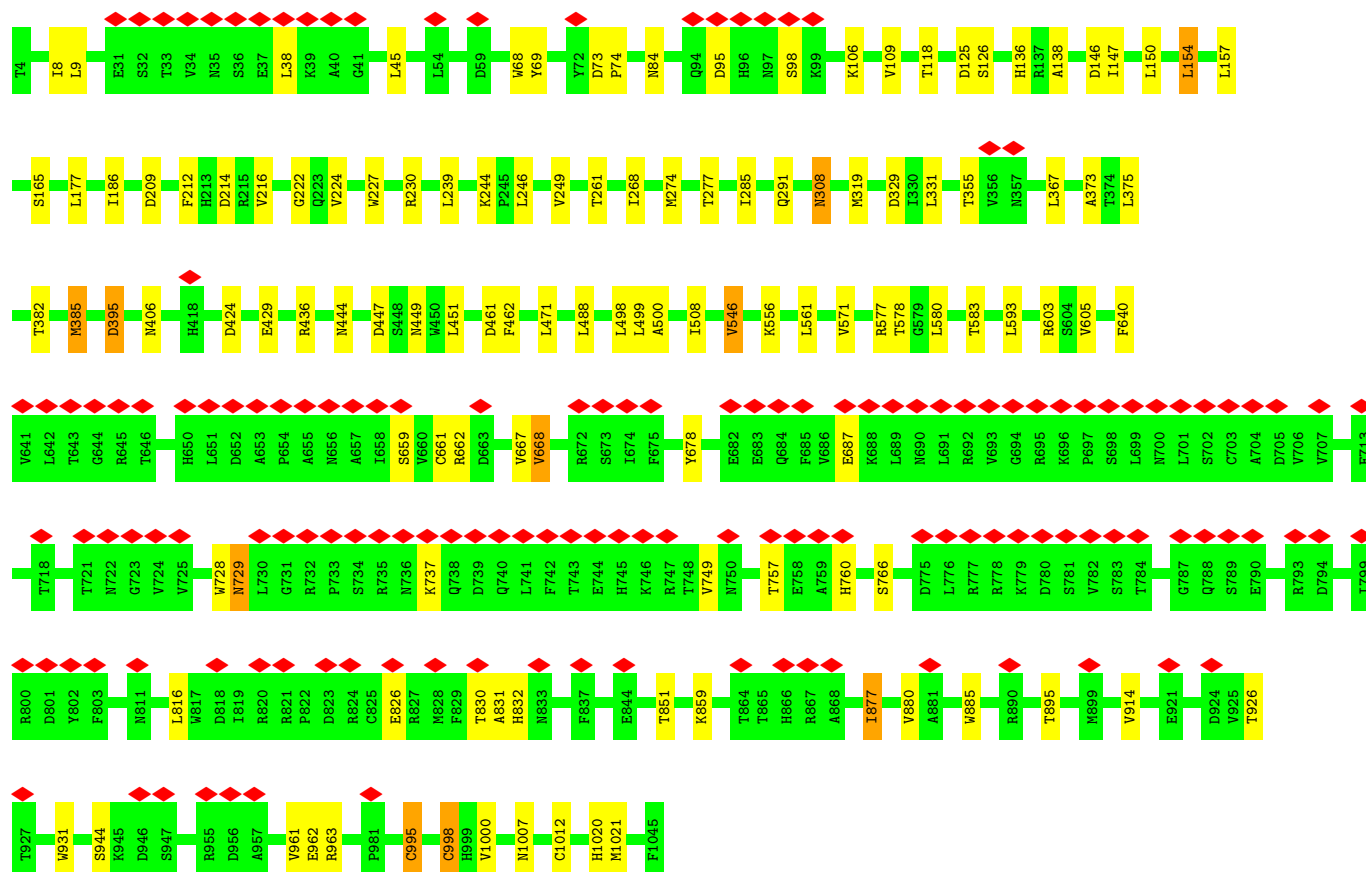
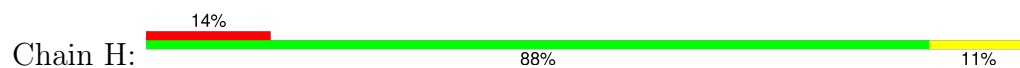
• Molecule 4: Cytosolic arginine sensor for mTORC1 subunit 1



• Molecule 4: Cytosolic arginine sensor for mTORC1 subunit 1



• Molecule 5: GATOR2 complex protein WDR24



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	865410	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 ($e^-/\text{\AA}^2$)	69.44	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2100	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.567	Depositor
Minimum map value	-0.202	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.015	Depositor
Recommended contour level	0.05	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

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.07	0/6033	0.22	0/8213
1	B	0.08	0/5715	0.24	0/7805
2	C	0.08	0/4477	0.24	0/6086
3	D	0.08	0/4108	0.24	0/5609
4	E	0.10	0/2538	0.29	0/3473
4	F	0.09	0/2331	0.27	0/3196
4	G	0.08	0/2133	0.23	0/2918
5	H	0.08	0/7752	0.26	0/10626
All	All	0.08	0/35087	0.25	0/47926

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
5	H	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
5	H	395	ASP	Peptide

5.2 Too-close contacts

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

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5894	0	5516	29	0
1	B	5582	0	5051	33	0
2	C	4367	0	4153	21	0
3	D	4010	0	3697	21	0
4	E	2477	0	2389	11	0
4	F	2279	0	2147	22	0
4	G	2084	0	1939	14	0
5	H	7559	0	6775	57	0
All	All	34252	0	31667	203	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 (203) 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:H:995:CYS:SG	5:H:998:CYS:N	2.56	0.72
1:B:1197:CYS:SG	1:B:1200:CYS:N	2.63	0.71
3:D:892:GLN:HA	3:D:920:VAL:HG13	1.74	0.69
3:D:847:ILE:HD11	3:D:871:VAL:HG11	1.75	0.69
1:B:1197:CYS:SG	1:B:1198:HIS:N	2.67	0.67
5:H:728:TRP:HE3	5:H:737:LYS:HB2	1.59	0.67
1:B:775:ASP:OD1	1:B:782:ASN:ND2	2.28	0.66
5:H:447:ASP:OD1	5:H:449:ASN:ND2	2.27	0.66
5:H:603:ARG:NH2	5:H:640:PHE:O	2.29	0.66
4:E:296:TYR:HB3	4:E:307:LEU:HB2	1.78	0.65
1:A:419:GLN:NE2	1:A:441:HIS:O	2.29	0.65
1:A:533:MET:SD	1:A:533:MET:N	2.70	0.65
5:H:885:TRP:O	5:H:931:TRP:NE1	2.29	0.65
1:B:1185:LYS:O	1:B:1189:GLN:NE2	2.29	0.64
1:B:996:ASP:O	1:B:1000:ASN:ND2	2.30	0.64
5:H:998:CYS:SG	5:H:1020:HIS:ND1	2.65	0.64
1:B:533:MET:N	1:B:533:MET:SD	2.73	0.62
1:A:978:GLU:O	1:A:986:ARG:NH1	2.31	0.62
4:F:203:ASP:OD1	4:G:36:ARG:NH1	2.33	0.62
5:H:385:MET:SD	5:H:385:MET:N	2.69	0.62
4:F:36:ARG:NH2	4:G:207:TYR:O	2.32	0.62
1:B:826:ILE:HG22	1:B:1021:LEU:HA	1.80	0.62
5:H:678:TYR:HA	5:H:687:GLU:HA	1.80	0.62
1:B:1193:TRP:HD1	1:B:1194:PHE:H	1.47	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1069:ARG:NH1	3:D:1192:ARG:O	2.32	0.61
2:C:696:ASN:HD21	2:C:700:THR:H	1.48	0.60
5:H:961:VAL:O	5:H:962:GLU:N	2.34	0.60
2:C:443:GLN:OE1	2:C:463:LYS:NZ	2.34	0.60
2:C:528:VAL:HG22	2:C:541:THR:HG22	1.82	0.60
1:A:955:ASN:HD22	1:A:958:LEU:HG	1.66	0.60
4:E:308:VAL:HG22	4:E:309:PRO:HD2	1.84	0.60
5:H:222:GLY:HA2	5:H:246:LEU:HD13	1.83	0.60
3:D:928:ILE:HG23	3:D:929:THR:HG23	1.84	0.60
5:H:73:ASP:HB3	5:H:74:PRO:HD3	1.84	0.59
4:E:101:VAL:HG11	4:E:123:ILE:HD13	1.86	0.58
2:C:585:ARG:NH1	2:C:1045:GLU:OE1	2.36	0.58
3:D:920:VAL:O	3:D:938:GLY:N	2.37	0.57
1:B:662:HIS:HD2	1:B:666:VAL:HG12	1.68	0.57
5:H:998:CYS:HB2	5:H:1000:VAL:HG22	1.85	0.57
5:H:661:CYS:SG	5:H:662:ARG:N	2.78	0.57
1:A:1055:LEU:HD13	1:A:1055:LEU:H	1.70	0.57
1:B:367:MET:HE1	1:B:392:ILE:HG13	1.85	0.56
1:B:367:MET:HE2	1:B:412:MET:HE1	1.87	0.56
4:F:298:ILE:HG23	4:F:305:HIS:HB2	1.85	0.56
1:A:981:VAL:O	1:A:986:ARG:NH1	2.38	0.56
5:H:224:VAL:HB	5:H:239:LEU:HD12	1.87	0.56
1:A:1157:LEU:HD12	1:A:1208:HIS:HB3	1.89	0.55
1:B:392:ILE:HG22	1:B:394:ALA:H	1.70	0.55
4:F:186:VAL:HG12	4:F:240:VAL:HG22	1.88	0.55
3:D:1188:ASP:HA	3:D:1191:LYS:HE3	1.89	0.55
4:F:296:TYR:HB3	4:F:307:LEU:HB2	1.89	0.55
5:H:106:LYS:HE2	5:H:150:LEU:HD11	1.88	0.55
1:B:522:GLY:HA2	1:B:549:SER:HA	1.88	0.55
4:E:242:ASP:OD1	4:E:242:ASP:N	2.40	0.54
5:H:580:LEU:HD11	5:H:593:LEU:HD23	1.89	0.54
1:A:416:SER:OG	1:A:418:ASP:OD1	2.25	0.54
2:C:526:THR:OG1	2:C:542:CYS:SG	2.65	0.54
4:F:75:THR:OG1	4:F:127:GLU:OE1	2.25	0.54
5:H:571:VAL:HG22	5:H:583:THR:HG22	1.90	0.54
5:H:749:VAL:HA	5:H:766:SER:HA	1.89	0.54
5:H:436:ARG:NH2	5:H:462:PHE:O	2.41	0.53
5:H:216:VAL:HG22	5:H:227:TRP:HB2	1.90	0.53
5:H:268:ILE:HG12	5:H:291:GLN:HG2	1.91	0.53
4:E:283:GLN:N	4:E:283:GLN:OE1	2.42	0.53
4:G:47:THR:OG1	4:G:50:ASP:O	2.25	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:682:THR:HG23	3:D:683:ARG:HG3	1.90	0.53
5:H:177:LEU:HD13	5:H:186:ILE:HG12	1.91	0.53
2:C:795:ARG:HB2	2:C:800:LEU:HD13	1.91	0.52
4:F:99:ARG:HB2	4:F:289:ALA:HB1	1.92	0.52
4:G:3:LEU:HB3	4:G:143:ILE:HA	1.90	0.52
4:F:78:VAL:HG22	4:F:124:LEU:HG	1.91	0.52
2:C:696:ASN:HD22	2:C:696:ASN:H	1.57	0.52
4:F:183:ARG:HE	4:F:262:LEU:HD13	1.75	0.51
5:H:274:MET:HA	5:H:285:ILE:HD12	1.91	0.51
5:H:926:THR:HA	5:H:944:SER:HA	1.91	0.51
5:H:831:ALA:O	5:H:859:LYS:NZ	2.44	0.50
1:B:627:THR:HG23	1:B:629:ASP:H	1.77	0.50
1:B:792:GLN:OE1	1:B:1015:ASN:ND2	2.44	0.50
1:B:1154:ARG:NH1	1:B:1161:ASN:OD1	2.44	0.50
3:D:929:THR:OG1	3:D:931:ASN:OD1	2.24	0.50
4:G:229:ALA:HB3	4:G:240:VAL:HG13	1.93	0.49
5:H:118:THR:HG21	5:H:165:SER:HA	1.93	0.49
1:B:796:LEU:HD22	1:B:1016:LEU:HD13	1.94	0.49
1:A:623:LEU:HD21	1:A:625:ILE:HD11	1.94	0.49
4:E:229:ALA:HB3	4:E:240:VAL:HG13	1.94	0.49
1:A:857:ASN:OD1	1:A:858:GLU:N	2.43	0.49
1:B:1212:TRP:CD1	1:B:1220:PRO:HG3	2.48	0.49
5:H:136:HIS:HD2	5:H:138:ALA:H	1.62	0.48
5:H:249:VAL:HG22	5:H:261:THR:HG22	1.95	0.48
1:B:936:ASP:OD1	1:B:936:ASP:N	2.46	0.48
4:F:105:LEU:HD22	4:F:110:VAL:HG11	1.95	0.48
5:H:214:ASP:OD2	5:H:230:ARG:NH1	2.45	0.48
5:H:488:LEU:HD23	5:H:500:ALA:HB2	1.95	0.48
1:B:662:HIS:CE1	1:B:688:ARG:HD3	2.49	0.47
2:C:651:THR:OG1	2:C:652:LYS:N	2.47	0.47
5:H:125:ASP:OD1	5:H:126:SER:N	2.40	0.47
5:H:816:LEU:O	5:H:826:GLU:N	2.47	0.47
1:A:1143:THR:O	1:A:1143:THR:OG1	2.33	0.47
5:H:577:ARG:HH11	5:H:580:LEU:HB2	1.79	0.47
1:A:1011:LYS:HA	1:A:1031:LEU:HD23	1.95	0.47
4:F:120:THR:OG1	4:F:121:ASP:N	2.47	0.47
4:G:78:VAL:HG22	4:G:124:LEU:HG	1.95	0.47
4:G:117:THR:HG21	4:G:233:ILE:HD12	1.97	0.47
5:H:832:HIS:HE1	5:H:851:THR:HG23	1.79	0.47
1:B:441:HIS:NE2	1:B:463:SER:OG	2.39	0.47
1:B:659:PHE:HZ	1:B:698:TRP:HB2	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:288:LEU:HD21	4:F:316:VAL:HG23	1.96	0.47
3:D:867:HIS:HD2	3:D:891:SER:HB2	1.79	0.46
5:H:329:ASP:N	5:H:329:ASP:OD1	2.48	0.46
4:G:54:MET:SD	4:G:54:MET:N	2.89	0.46
5:H:546:VAL:HG23	5:H:561:LEU:HB2	1.98	0.46
2:C:1040:ASP:OD1	2:C:1040:ASP:N	2.47	0.46
4:E:181:GLN:HB2	4:E:223:SER:OG	2.15	0.46
5:H:961:VAL:O	5:H:963:ARG:N	2.45	0.46
3:D:726:TYR:HD2	3:D:774:LEU:HD13	1.81	0.46
5:H:68:TRP:HD1	5:H:69:TYR:H	1.64	0.46
1:B:775:ASP:OD1	1:B:775:ASP:N	2.50	0.45
3:D:939:ASP:N	3:D:939:ASP:OD1	2.49	0.45
1:A:627:THR:HG21	1:A:631:ARG:HE	1.82	0.45
3:D:652:GLY:HA2	3:D:996:ALA:HA	1.98	0.45
1:B:857:ASN:OD1	1:B:858:GLU:N	2.46	0.45
2:C:650:ALA:HB2	2:C:693:VAL:HG13	1.98	0.45
5:H:95:ASP:O	5:H:98:SER:OG	2.33	0.45
2:C:696:ASN:ND2	2:C:701:VAL:H	2.15	0.45
4:F:300:THR:OG1	4:F:301:PHE:N	2.50	0.45
5:H:729:ASN:H	5:H:737:LYS:HB3	1.82	0.45
1:A:552:CYS:SG	1:A:610:ILE:HG12	2.57	0.44
4:G:95:THR:O	4:G:99:ARG:N	2.43	0.44
5:H:146:ASP:OD1	5:H:147:ILE:N	2.49	0.44
1:B:542:HIS:CE1	1:B:593:LYS:HA	2.53	0.44
4:F:173:THR:HA	4:F:283:GLN:HE22	1.83	0.44
5:H:880:VAL:HG13	5:H:895:THR:HG23	2.00	0.44
4:F:190:ASP:OD1	4:F:193:THR:OG1	2.33	0.44
3:D:539:THR:HG21	3:D:555:ARG:HD3	1.99	0.44
3:D:921:VAL:HA	3:D:937:GLY:HA2	2.00	0.43
1:A:1112:LYS:HD2	1:A:1147:GLY:HA3	2.00	0.43
4:F:190:ASP:OD1	4:F:190:ASP:N	2.34	0.43
5:H:395:ASP:N	5:H:395:ASP:OD1	2.51	0.43
1:A:466:ARG:NH1	1:A:498:ARG:O	2.39	0.43
1:A:1210:LEU:HD11	2:C:1141:MET:HE2	2.00	0.43
4:F:264:ARG:HE	4:F:313:ILE:HD12	1.84	0.43
5:H:308:ASN:OD1	5:H:308:ASN:N	2.51	0.43
5:H:659:SER:HB3	5:H:668:VAL:HG13	2.01	0.43
1:A:627:THR:HG23	1:A:629:ASP:H	1.83	0.43
1:A:669:VAL:HG13	1:A:678:LEU:HD11	2.00	0.43
1:B:1209:MET:HE3	1:B:1213:PHE:HE2	1.84	0.43
4:F:13:LEU:HD22	4:F:71:VAL:HG12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:105:LEU:HB3	4:F:110:VAL:HB	2.00	0.43
1:A:1110:CYS:SG	1:A:1112:LYS:NZ	2.85	0.43
2:C:1143:TRP:CD1	2:C:1151:PRO:HG3	2.54	0.43
4:G:94:VAL:HG23	4:G:96:LYS:H	1.84	0.43
5:H:444:ASN:HB3	5:H:447:ASP:O	2.19	0.43
2:C:696:ASN:HD21	2:C:700:THR:N	2.13	0.43
4:E:304:ASP:N	4:E:304:ASP:OD1	2.51	0.42
5:H:1007:ASN:OD1	5:H:1007:ASN:N	2.50	0.42
3:D:746:ASN:OD1	3:D:746:ASN:N	2.52	0.42
2:C:1090:CYS:SG	2:C:1091:SER:N	2.92	0.42
5:H:146:ASP:HB2	5:H:154:LEU:HD23	2.00	0.42
5:H:367:LEU:HD11	5:H:373:ALA:HB2	2.01	0.42
3:D:843:CYS:HB3	3:D:870:TRP:CZ3	2.55	0.42
2:C:800:LEU:HD23	2:C:1030:VAL:HG13	2.01	0.42
5:H:68:TRP:HD1	5:H:69:TYR:N	2.18	0.42
1:A:976:LEU:HB3	1:A:1002:TYR:CE1	2.54	0.42
1:A:1140:SER:OG	1:A:1141:LYS:N	2.53	0.42
3:D:811:ALA:HB2	3:D:873:ASP:HA	2.01	0.42
3:D:788:LEU:HB3	3:D:796:TRP:CE3	2.55	0.42
4:E:120:THR:OG1	4:E:121:ASP:N	2.52	0.42
3:D:845:ASN:OD1	3:D:845:ASN:N	2.53	0.42
4:G:6:LEU:HD23	4:G:8:HIS:NE2	2.35	0.42
5:H:84:ASN:OD1	5:H:84:ASN:N	2.48	0.42
1:A:485:GLN:O	1:A:487:HIS:ND1	2.50	0.41
1:A:883:ASN:O	1:A:886:VAL:HG12	2.20	0.41
4:F:29:LYS:HG3	4:F:36:ARG:HG3	2.02	0.41
5:H:499:LEU:HD13	5:H:508:ILE:HG12	2.02	0.41
4:F:256:THR:OG1	4:F:257:SER:N	2.51	0.41
1:A:935:THR:HG22	1:A:937:GLU:H	1.85	0.41
3:D:995:ILE:HB	3:D:996:ALA:H	1.62	0.41
5:H:914:VAL:HG11	5:H:961:VAL:HG13	2.02	0.41
1:B:806:TYR:CD1	1:B:822:VAL:HG21	2.55	0.41
1:B:1193:TRP:CD1	1:B:1194:PHE:H	2.33	0.41
4:E:6:LEU:HB2	4:E:78:VAL:HG13	2.03	0.41
4:G:103:ALA:HB3	4:G:104:PRO:HD3	2.02	0.41
1:A:1102:GLN:HE22	2:C:1134:GLY:H	1.69	0.41
1:B:672:ASN:HD21	1:B:674:THR:HG22	1.86	0.41
2:C:580:ASN:HA	2:C:581:PRO:HD3	1.91	0.41
1:B:1035:TYR:O	1:B:1039:THR:HG22	2.21	0.41
4:G:287:PRO:HB3	4:G:319:VAL:HG12	2.03	0.41
5:H:274:MET:SD	5:H:274:MET:N	2.93	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:LYS:HE2	1:A:490:LYS:HB3	1.92	0.41
1:A:1016:LEU:HD12	1:A:1016:LEU:HA	1.95	0.41
2:C:358:SER:OG	2:C:391:GLY:O	2.34	0.41
1:B:452:HIS:CG	1:B:453:PRO:HD2	2.56	0.40
4:E:243:ALA:O	4:E:246:GLN:HG2	2.21	0.40
4:G:10:VAL:HG13	4:G:76:TRP:HB2	2.02	0.40
3:D:767:TRP:CZ3	3:D:776:LEU:HB2	2.56	0.40
5:H:209:ASP:HB3	5:H:212:PHE:O	2.22	0.40
5:H:877:ILE:HD13	5:H:877:ILE:H	1.86	0.40
2:C:983:SER:OG	2:C:984:SER:N	2.54	0.40
5:H:429:GLU:HG2	5:H:471:LEU:HD23	2.03	0.40
1:A:1201:ARG:NH2	2:C:1098:SER:O	2.54	0.40
1:B:899:LEU:HD22	1:B:961:MET:HG3	2.04	0.40
4:F:102:ILE:HD11	4:F:114:MET:HE1	2.04	0.40
1:B:690:TRP:CZ3	1:B:700:CYS:HB3	2.56	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	774/786 (98%)	746 (96%)	28 (4%)	0	100	100
1	B	747/786 (95%)	698 (93%)	49 (7%)	0	100	100
2	C	558/570 (98%)	542 (97%)	16 (3%)	0	100	100
3	D	536/545 (98%)	512 (96%)	24 (4%)	0	100	100
4	E	326/328 (99%)	309 (95%)	17 (5%)	0	100	100
4	F	311/328 (95%)	295 (95%)	16 (5%)	0	100	100
4	G	276/328 (84%)	271 (98%)	5 (2%)	0	100	100
5	H	1020/1042 (98%)	970 (95%)	50 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	4548/4713 (96%)	4343 (96%)	205 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	593/684 (87%)	580 (98%)	13 (2%)	45	63
1	B	538/684 (79%)	514 (96%)	24 (4%)	24	51
2	C	466/492 (95%)	447 (96%)	19 (4%)	27	52
3	D	387/462 (84%)	379 (98%)	8 (2%)	47	64
4	E	260/285 (91%)	249 (96%)	11 (4%)	26	52
4	F	221/285 (78%)	213 (96%)	8 (4%)	31	56
4	G	200/285 (70%)	194 (97%)	6 (3%)	36	59
5	H	709/923 (77%)	673 (95%)	36 (5%)	21	48
All	All	3374/4100 (82%)	3249 (96%)	125 (4%)	31	55

All (125) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	378	THR
1	A	475	VAL
1	A	504	VAL
1	A	533	MET
1	A	623	LEU
1	A	699	LYS
1	A	862	LEU
1	A	1019	ILE
1	A	1055	LEU
1	A	1058	LEU
1	A	1064	GLN
1	A	1108	ASN

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Mol	Chain	Res	Type
1	A	1209	MET
1	B	358	LEU
1	B	400	ILE
1	B	462	CYS
1	B	470	VAL
1	B	518	CYS
1	B	533	MET
1	B	544	ILE
1	B	603	VAL
1	B	616	LEU
1	B	653	ILE
1	B	662	HIS
1	B	666	VAL
1	B	701	THR
1	B	769	LEU
1	B	834	LEU
1	B	842	HIS
1	B	876	VAL
1	B	909	GLN
1	B	912	ASN
1	B	931	LEU
1	B	1025	THR
1	B	1148	CYS
1	B	1197	CYS
1	B	1200	CYS
2	C	447	VAL
2	C	516	THR
2	C	557	MET
2	C	559	LEU
2	C	561	GLN
2	C	591	ILE
2	C	617	LYS
2	C	634	ILE
2	C	639	ASN
2	C	654	VAL
2	C	696	ASN
2	C	715	LYS
2	C	790	TYR
2	C	1042	GLN
2	C	1054	LEU
2	C	1103	VAL
2	C	1109	ARG

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Mol	Chain	Res	Type
2	C	1121	VAL
2	C	1126	VAL
3	D	550	LEU
3	D	576	LEU
3	D	649	ILE
3	D	668	HIS
3	D	719	VAL
3	D	836	LYS
3	D	845	ASN
3	D	889	SER
4	E	30	LEU
4	E	130	LEU
4	E	152	VAL
4	E	154	VAL
4	E	175	HIS
4	E	240	VAL
4	E	242	ASP
4	E	261	GLU
4	E	268	ILE
4	E	281	VAL
4	E	308	VAL
4	F	69	LEU
4	F	79	LEU
4	F	81	VAL
4	F	190	ASP
4	F	201	LEU
4	F	226	THR
4	F	268	ILE
4	F	313	ILE
4	G	17	ARG
4	G	107	GLU
4	G	115	LEU
4	G	137	LEU
4	G	145	ARG
4	G	240	VAL
5	H	8	ILE
5	H	9	LEU
5	H	38	LEU
5	H	45	LEU
5	H	109	VAL
5	H	154	LEU
5	H	157	LEU

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Mol	Chain	Res	Type
5	H	244	LYS
5	H	277	THR
5	H	308	ASN
5	H	319	MET
5	H	331	LEU
5	H	355	THR
5	H	375	LEU
5	H	382	THR
5	H	385	MET
5	H	406	ASN
5	H	424	ASP
5	H	451	LEU
5	H	461	ASP
5	H	498	LEU
5	H	546	VAL
5	H	556	LYS
5	H	578	THR
5	H	605	VAL
5	H	667	VAL
5	H	668	VAL
5	H	729	ASN
5	H	757	THR
5	H	760	HIS
5	H	830	THR
5	H	877	ILE
5	H	995	CYS
5	H	998	CYS
5	H	1012	CYS
5	H	1021	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (76) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	396	HIS
1	A	419	GLN
1	A	457	GLN
1	A	485	GLN
1	A	556	ASN
1	A	563	HIS
1	A	583	GLN
1	A	615	ASN
1	A	658	GLN

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Mol	Chain	Res	Type
1	A	662	HIS
1	A	663	ASN
1	A	697	ASN
1	A	783	HIS
1	A	883	ASN
1	A	887	GLN
1	A	952	GLN
1	A	955	ASN
1	A	998	GLN
1	A	1064	GLN
1	A	1102	GLN
1	A	1191	ASN
1	A	1192	ASN
1	A	1230	GLN
1	B	457	GLN
1	B	541	GLN
1	B	542	HIS
1	B	556	ASN
1	B	658	GLN
1	B	661	ASN
1	B	662	HIS
1	B	665	GLN
1	B	855	ASN
1	B	1000	ASN
1	B	1015	ASN
1	B	1043	GLN
1	B	1051	GLN
1	B	1064	GLN
2	C	432	HIS
2	C	481	GLN
2	C	511	HIS
2	C	686	HIS
2	C	696	ASN
2	C	818	GLN
2	C	1018	GLN
2	C	1022	GLN
2	C	1044	GLN
2	C	1129	GLN
2	C	1138	GLN
2	C	1156	HIS
3	D	662	ASN
3	D	805	HIS

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Mol	Chain	Res	Type
3	D	810	ASN
3	D	923	HIS
3	D	1012	ASN
3	D	1078	HIS
4	E	161	ASN
4	E	175	HIS
4	F	175	HIS
4	F	181	GLN
4	F	182	ASN
4	F	283	GLN
4	G	246	GLN
5	H	13	HIS
5	H	29	HIS
5	H	316	ASN
5	H	379	ASN
5	H	393	ASN
5	H	449	ASN
5	H	589	ASN
5	H	606	GLN
5	H	623	ASN
5	H	656	ASN
5	H	750	ASN
5	H	832	HIS
5	H	999	HIS
5	H	1028	GLN

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
5	H	10
2	C	5
1	A	5
3	D	4

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	640:PHE	C	641:VAL	N	114.11
1	H	325:PHE	C	326:THR	N	99.87
1	H	979:PHE	C	980:PRO	N	67.54
1	D	1099:SER	C	1174:LEU	N	44.51
1	D	1174:LEU	C	1178:ASP	N	39.20
1	C	830:PRO	C	982:SER	N	38.98
1	C	730:GLY	C	778:GLY	N	23.32
1	A	1166:VAL	C	1182:SER	N	22.99
1	A	711:VAL	C	757:GLU	N	22.33
1	H	981:PRO	C	982:PHE	N	19.06
1	C	361:ALA	C	387:GLU	N	16.86
1	H	470:CYS	C	471:LEU	N	14.57
1	H	148:CYS	C	149:LEU	N	14.37
1	A	1116:TYR	C	1140:SER	N	14.17
1	D	581:THR	C	630:LYS	N	12.84
1	A	846:GLY	C	852:ASP	N	11.82
1	H	991:SER	C	992:THR	N	11.02
1	D	957:SER	C	995:ILE	N	9.98
1	H	282:GLY	C	283:PRO	N	9.26
1	H	598:HIS	C	599:THR	N	9.01
1	C	404:PRO	C	410:MET	N	6.53
1	A	379:GLU	C	386:MET	N	4.81

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	499:VAL	C	510:SER	N	4.15
1	H	961:VAL	C	962:GLU	N	3.32

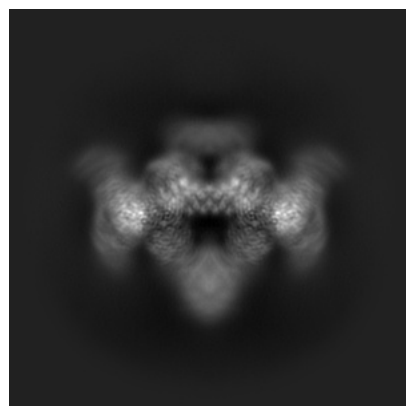
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-47278. These allow visual inspection of the internal detail of the map and identification of artifacts.

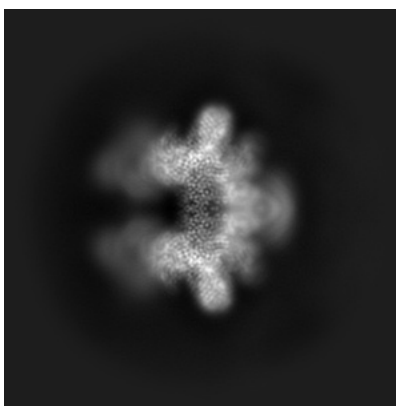
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

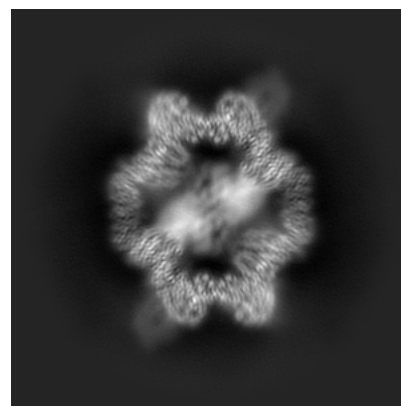
6.1.1 Primary map



X

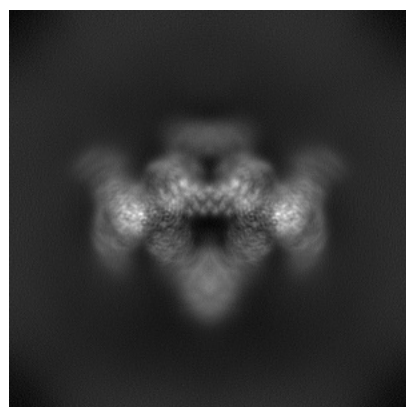


Y

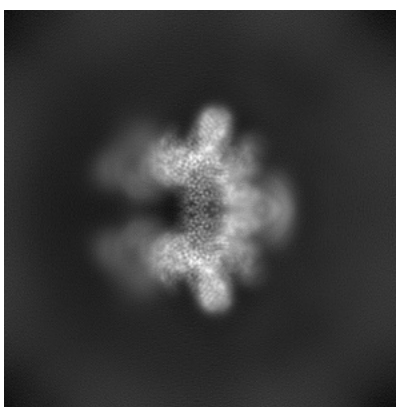


Z

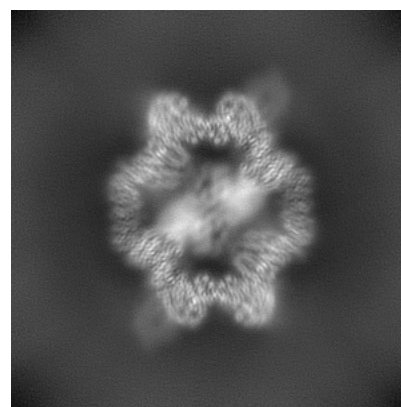
6.1.2 Raw map



X



Y

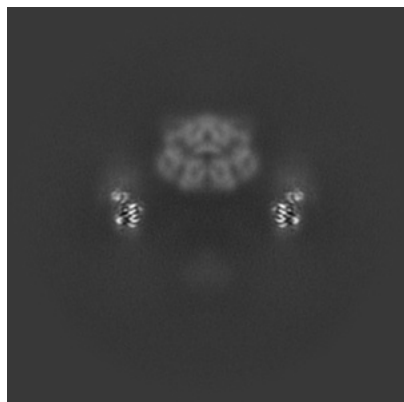


Z

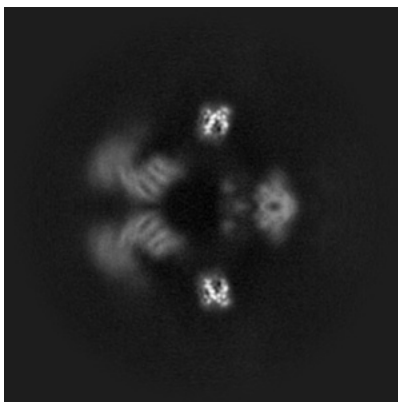
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

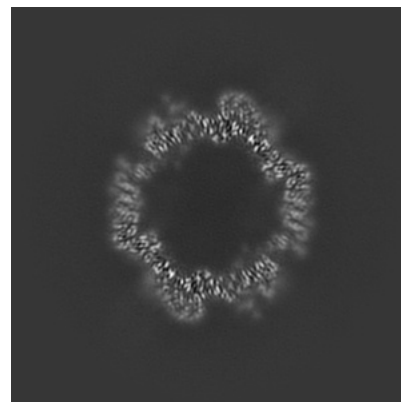
6.2.1 Primary map



X Index: 240

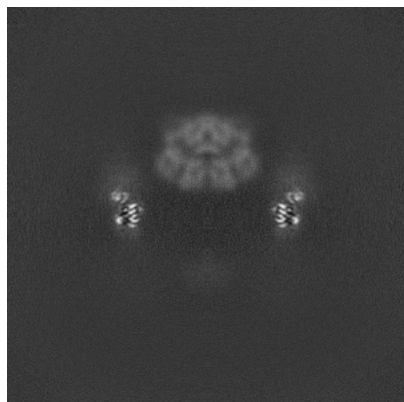


Y Index: 240

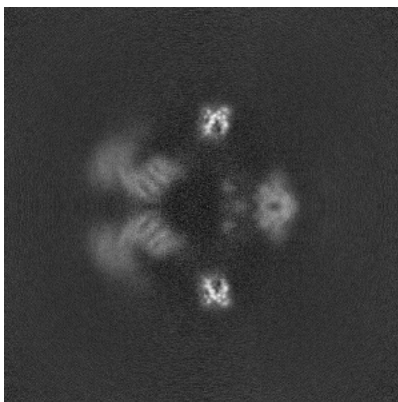


Z Index: 240

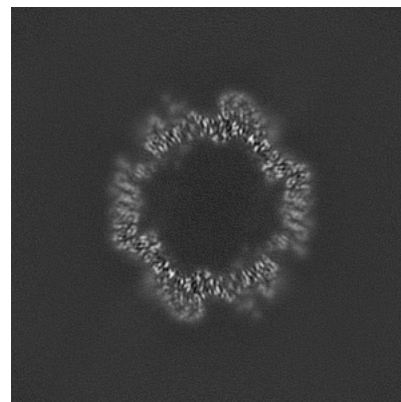
6.2.2 Raw map



X Index: 240



Y Index: 240

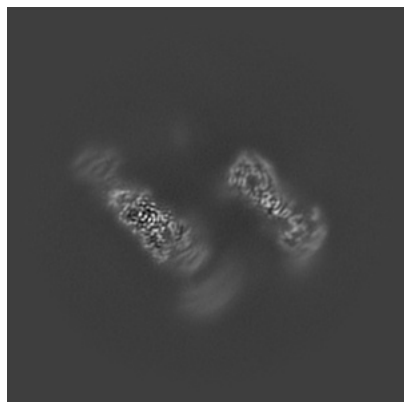


Z Index: 240

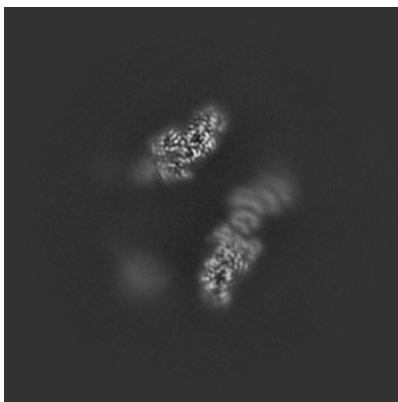
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

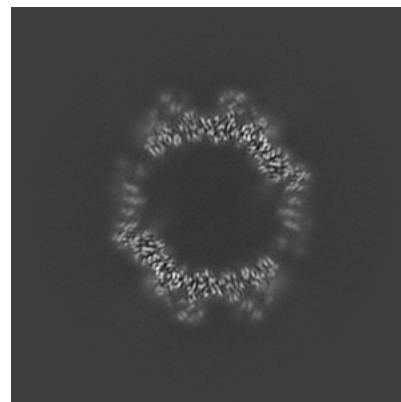
6.3.1 Primary map



X Index: 184

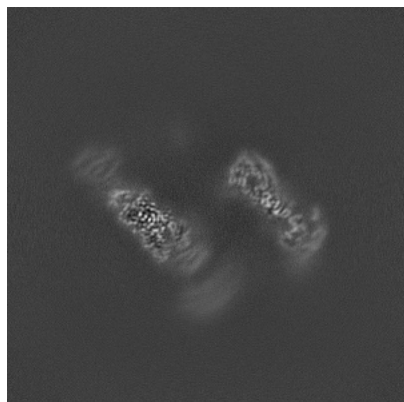


Y Index: 283

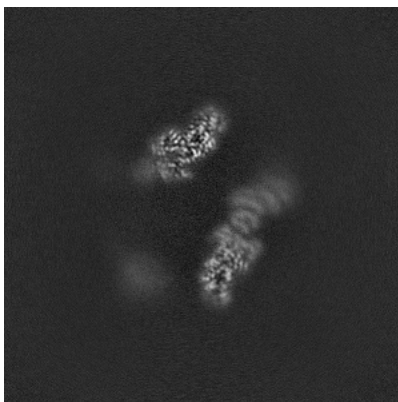


Z Index: 235

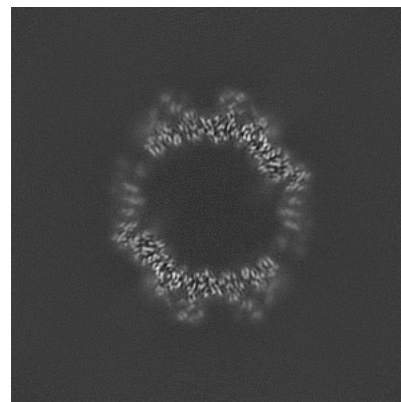
6.3.2 Raw map



X Index: 184



Y Index: 283

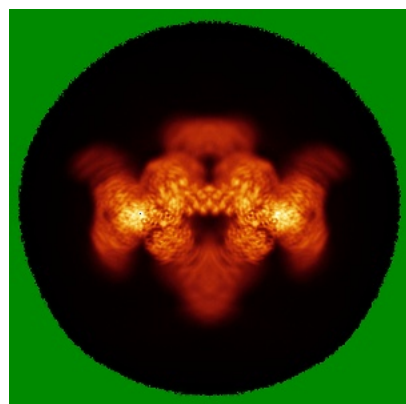


Z Index: 235

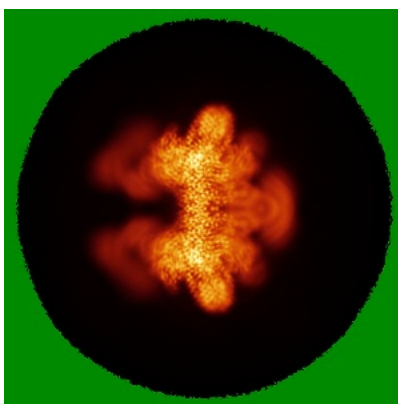
The images above show the largest variance slices of the map in three orthogonal directions.

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

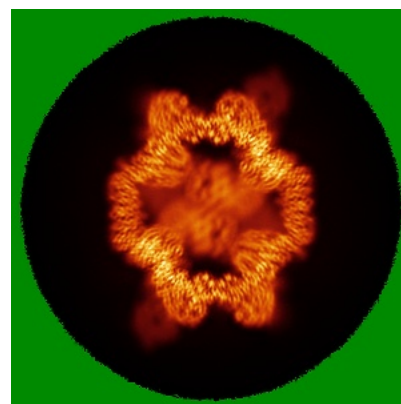
6.4.1 Primary map



X

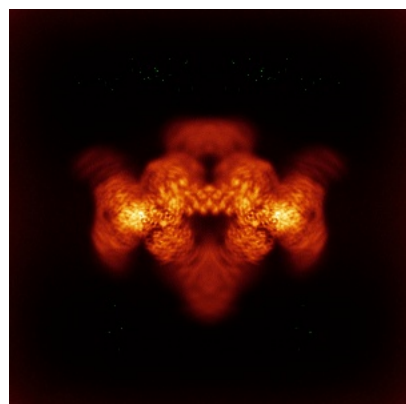


Y

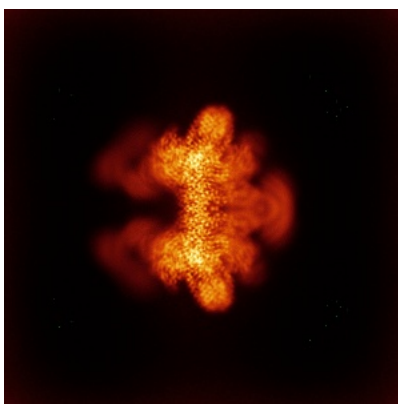


Z

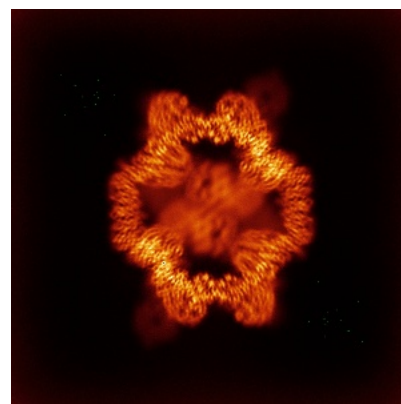
6.4.2 Raw map



X



Y

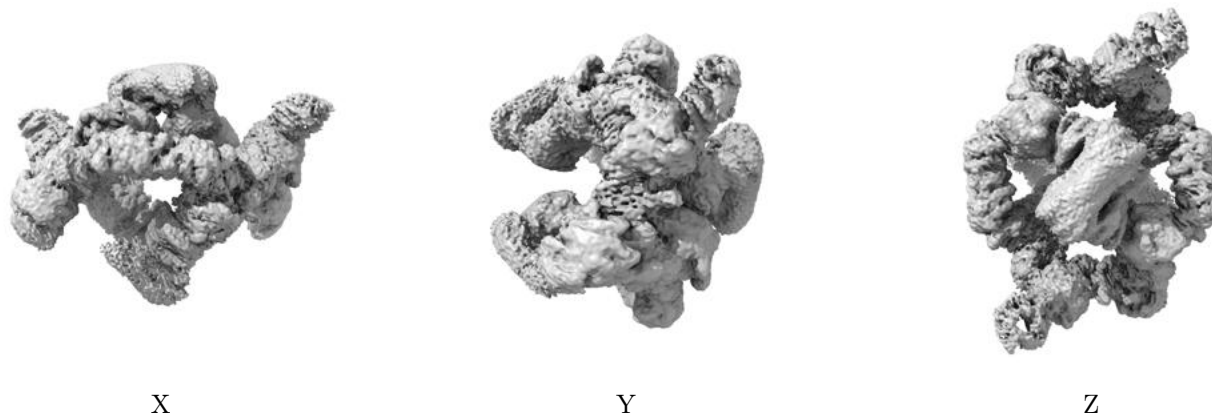


Z

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.

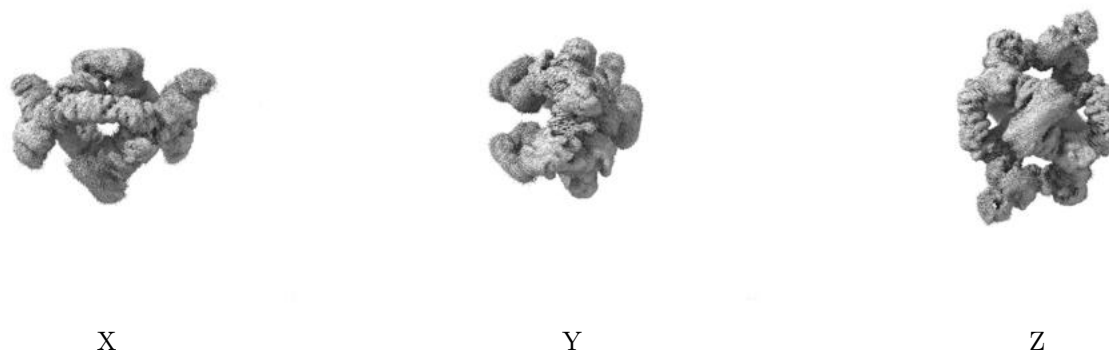
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.05. 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.

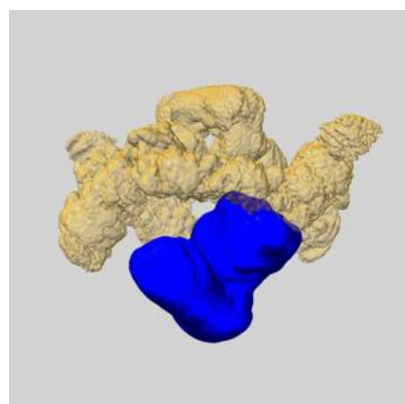
6.6 Mask visualisation [i](#)

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

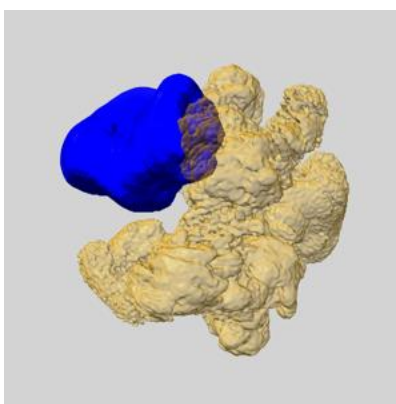
A mask typically either:

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

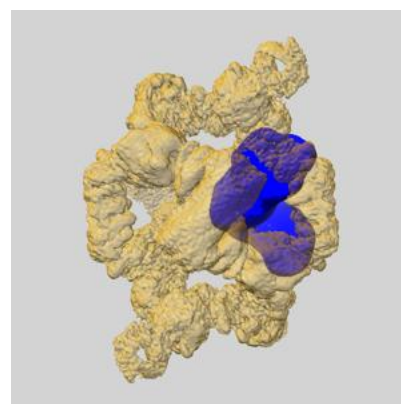
6.6.1 emd_47278_msk_5.map [i](#)



X

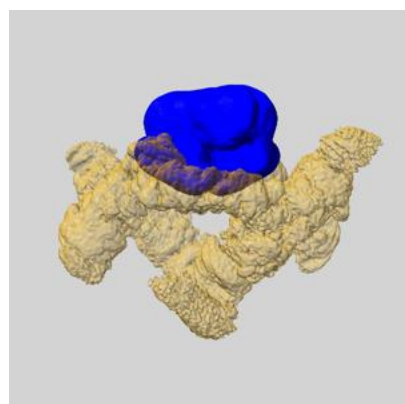


Y

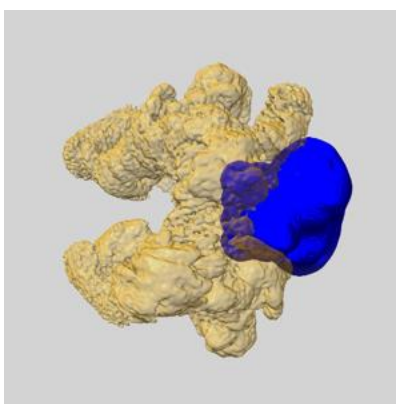


Z

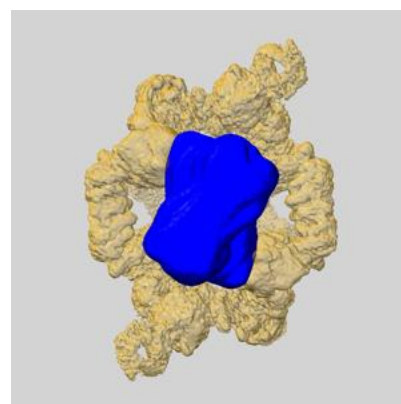
6.6.2 emd_47278_msk_4.map [i](#)



X

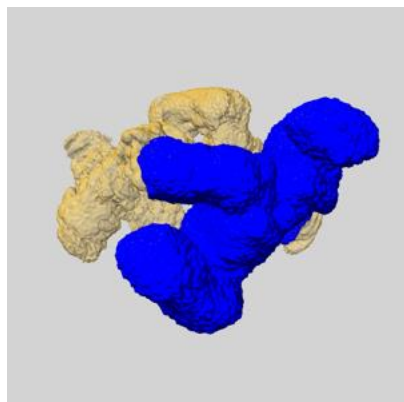


Y

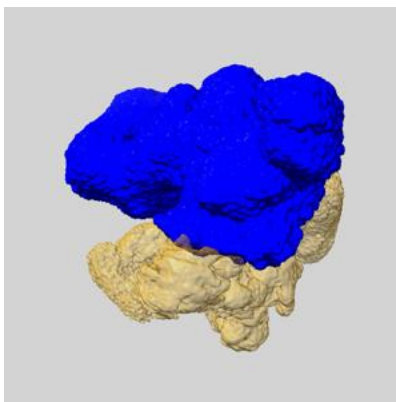


Z

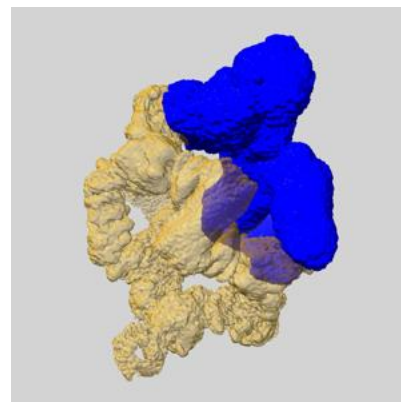
6.6.3 emd_47278_msk_3.map [i](#)



X

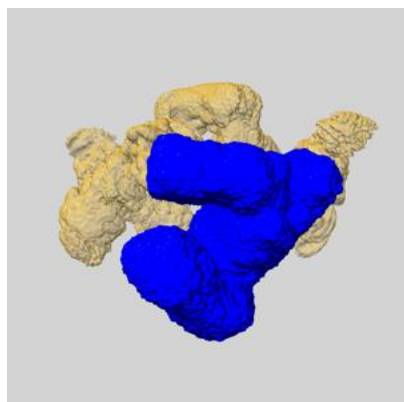


Y

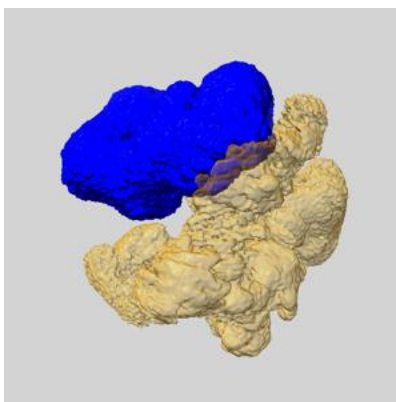


Z

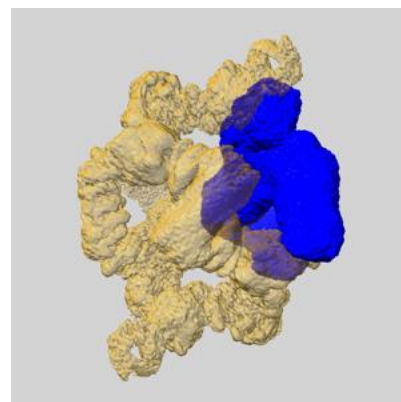
6.6.4 emd_47278_msk_2.map [i](#)



X

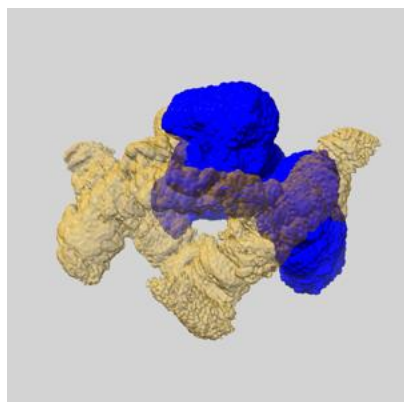


Y

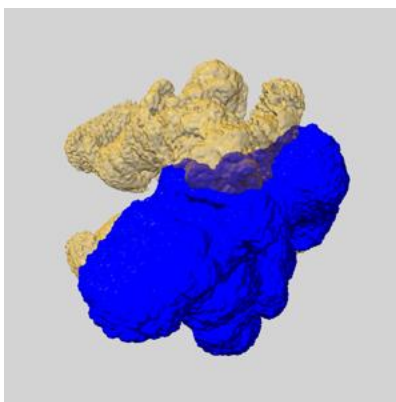


Z

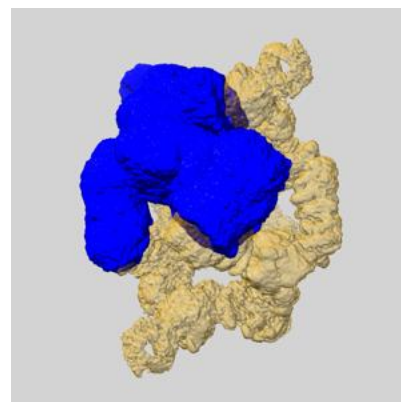
6.6.5 emd_47278_msk_1.map [i](#)



X



Y

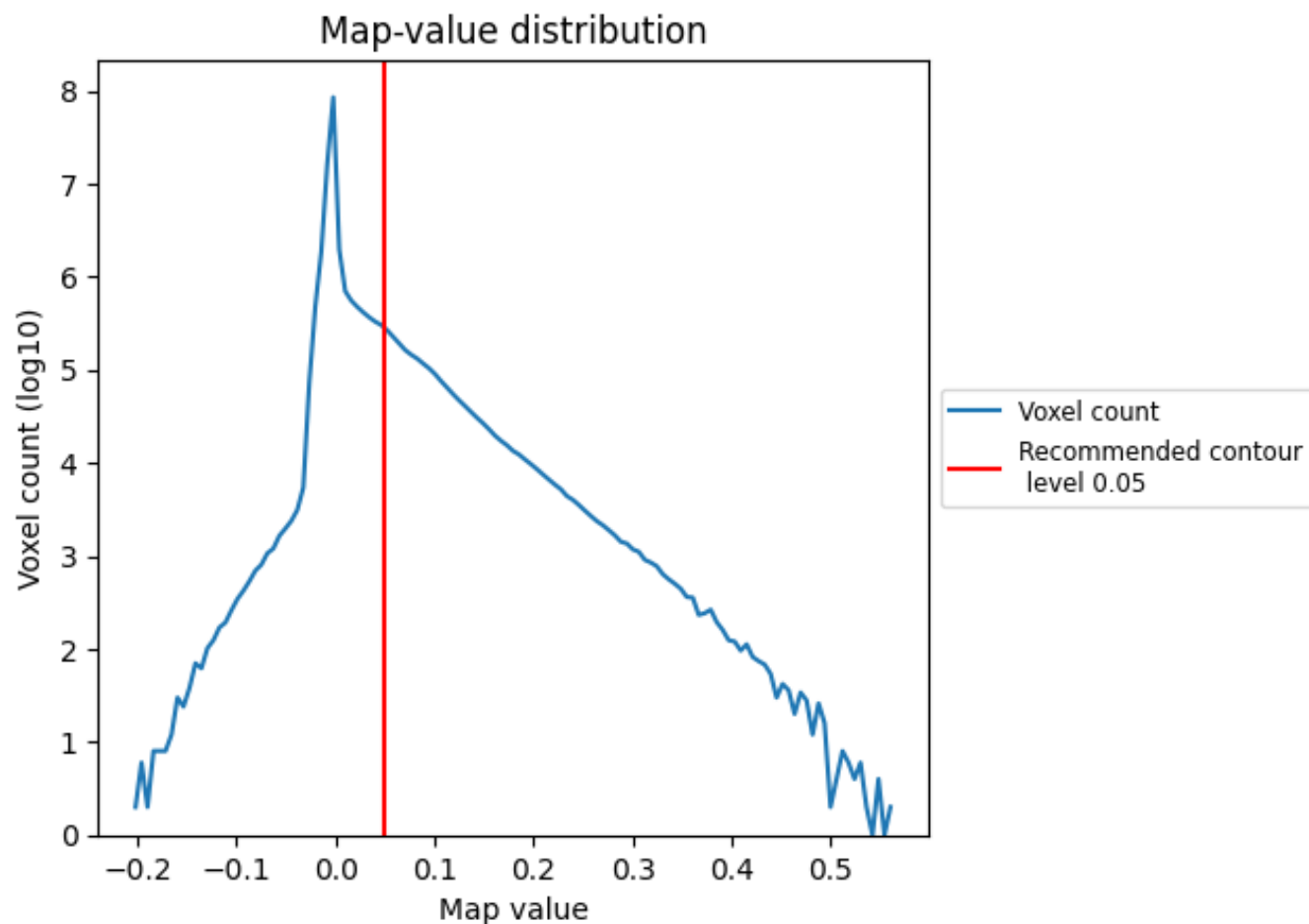


Z

7 Map analysis [i](#)

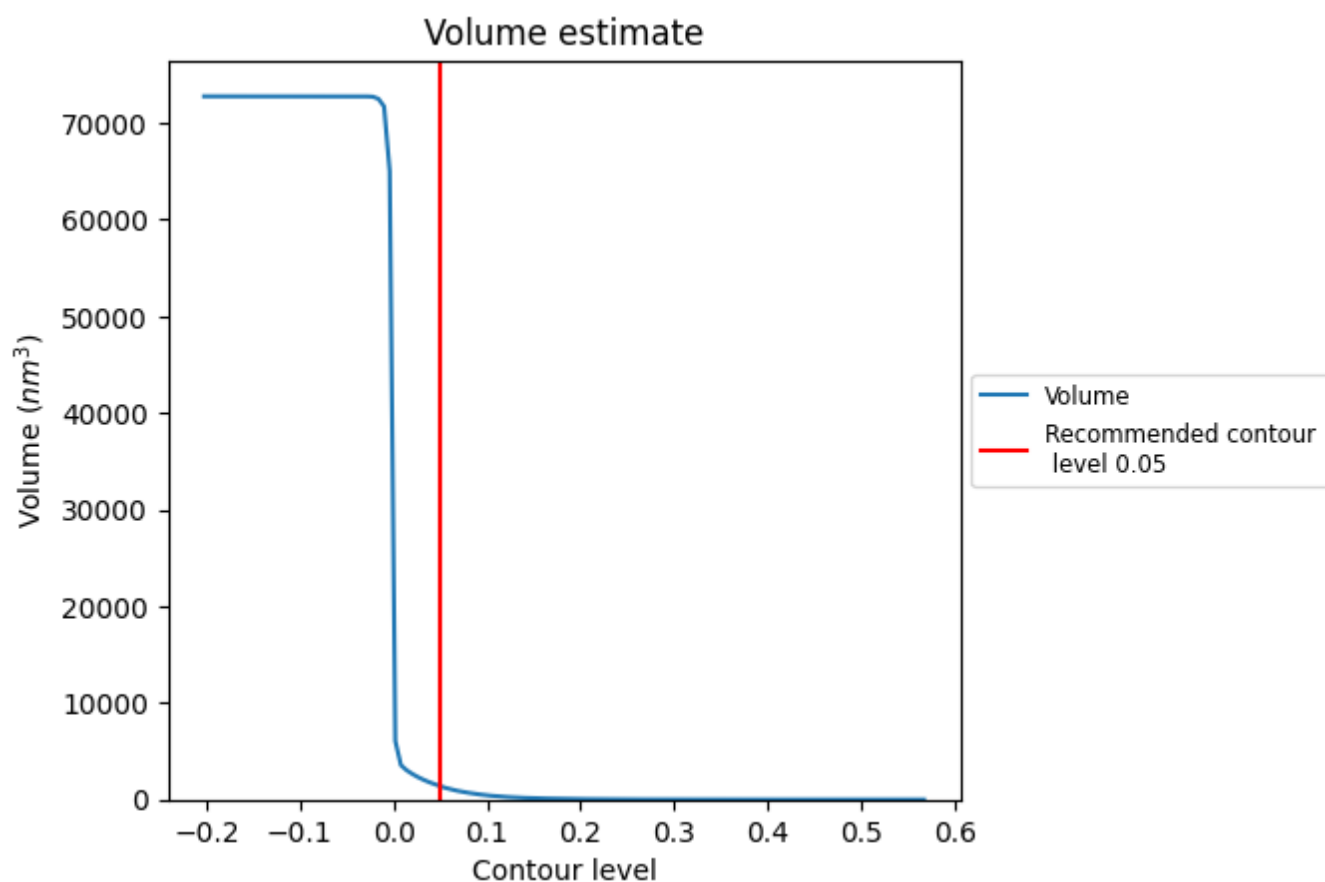
This section contains the results of statistical analysis of the map.

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.

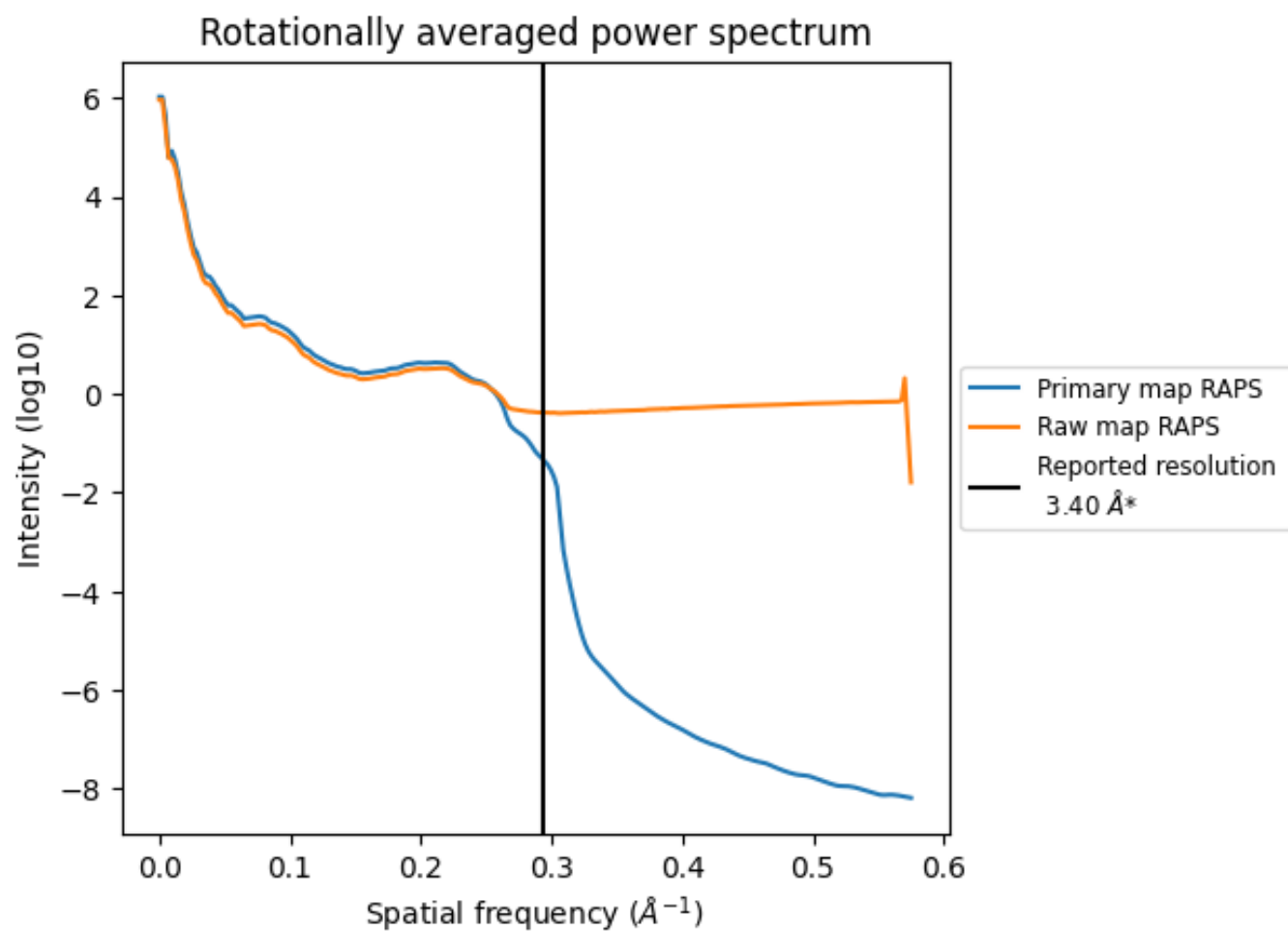
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1385 nm³; this corresponds to an approximate mass of 1251 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 ⓘ

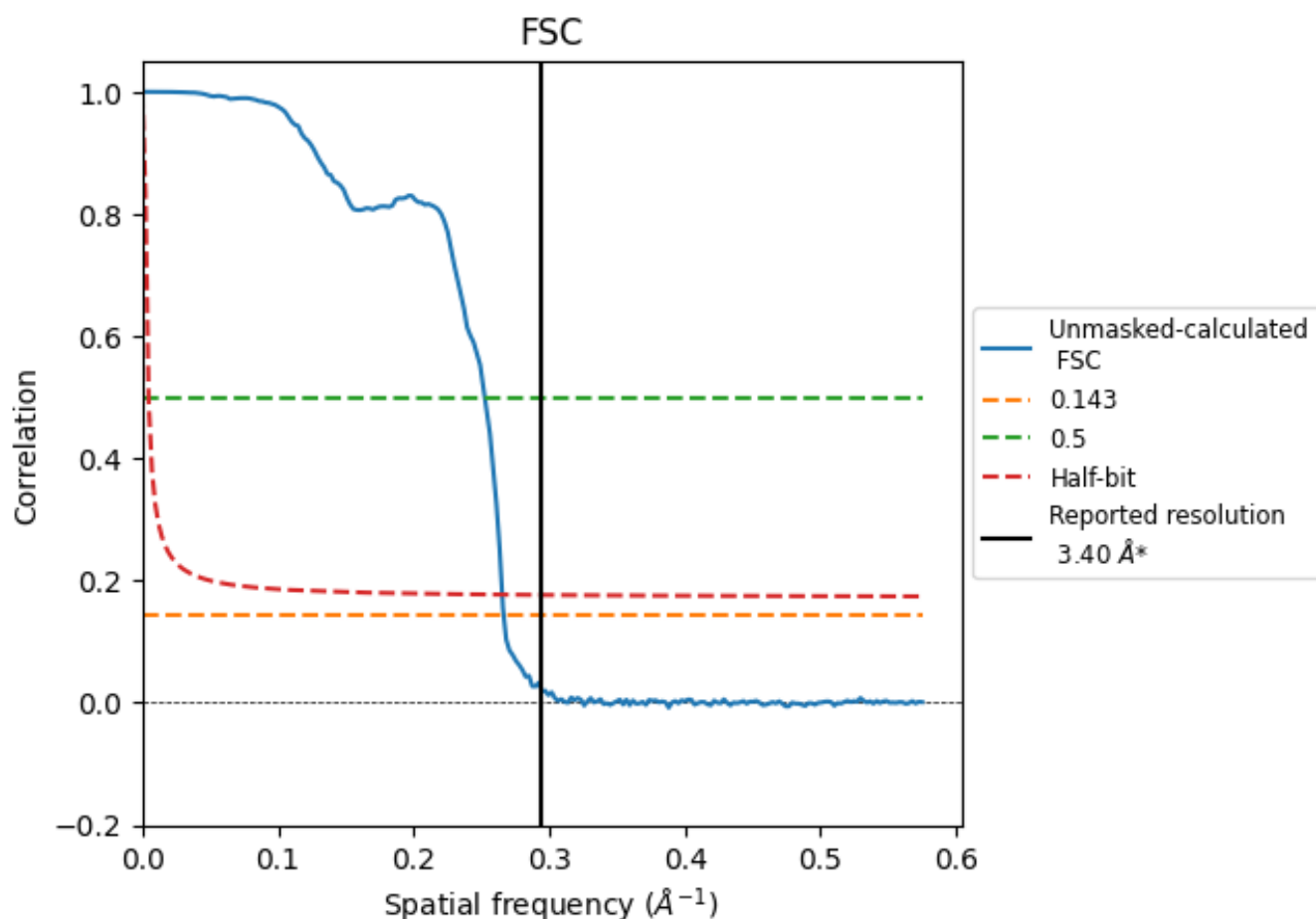


*Reported resolution corresponds to spatial frequency of 0.294 Å⁻¹

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 0.294 \AA^{-1}

8.2 Resolution estimates [i](#)

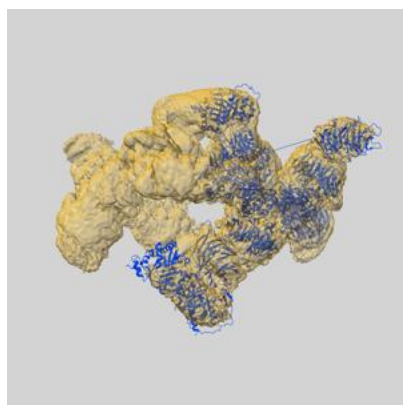
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.75	3.96	3.77

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.75 differs from the reported value 3.4 by more than 10 %

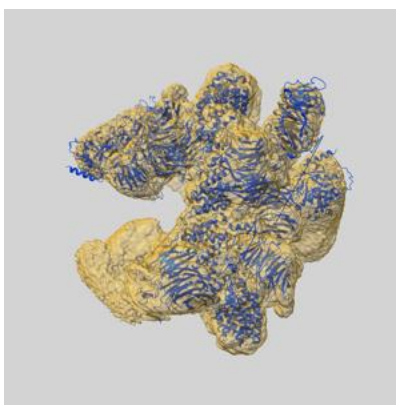
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-47278 and PDB model 9DX2. Per-residue inclusion information can be found in section [3](#) on page [5](#).

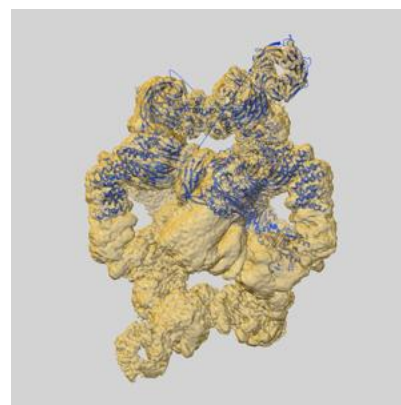
9.1 Map-model overlay [i](#)



X



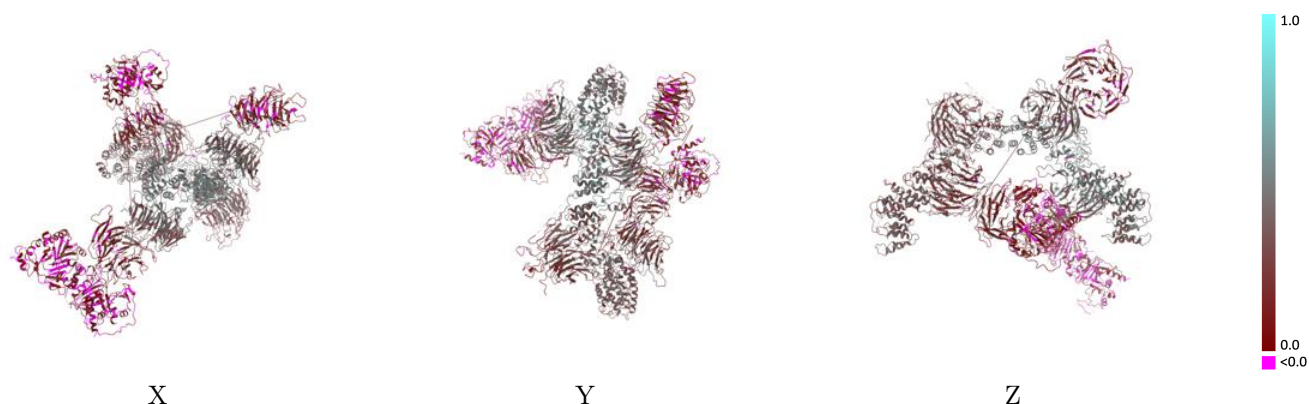
Y



Z

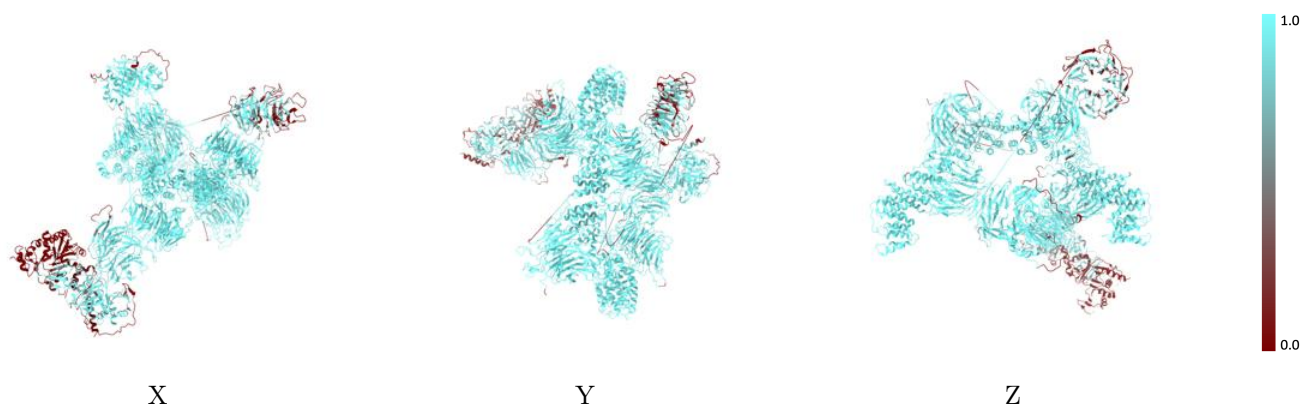
The images above show the 3D surface view of the map at the recommended contour level 0.05 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



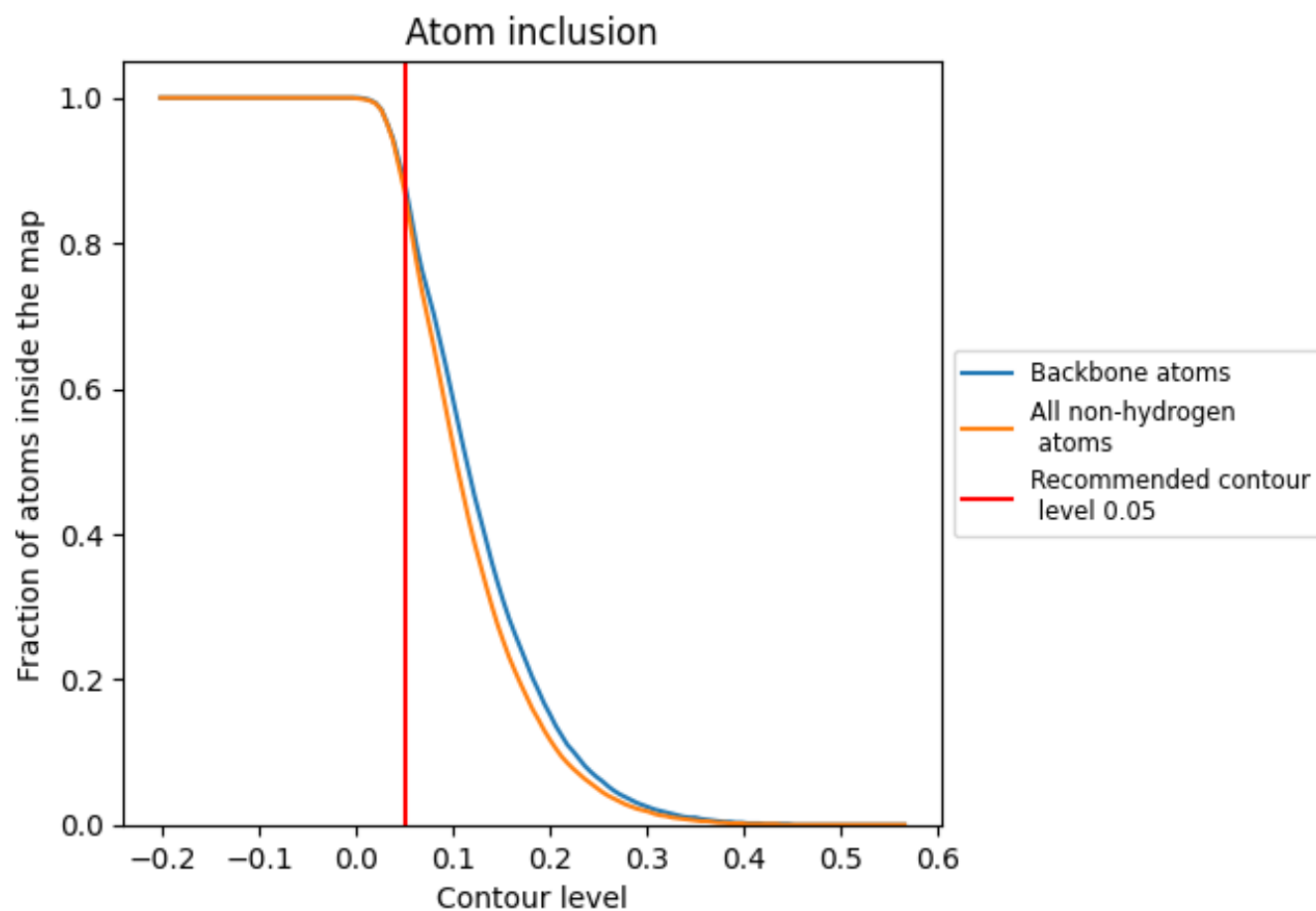
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.05).

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 87% 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 (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.8720	<div></div> 0.2770
A	<div></div> 0.9800	<div></div> 0.4250
B	<div></div> 0.9840	<div></div> 0.3380
C	<div></div> 0.9750	<div></div> 0.4320
D	<div></div> 0.9670	<div></div> 0.3190
E	<div></div> 0.8750	<div></div> 0.0940
F	<div></div> 0.6320	<div></div> 0.0540
G	<div></div> 0.2270	<div></div> 0.0570
H	<div></div> 0.8450	<div></div> 0.1940

1.0

0.0

<0.0