



## Full wwPDB EM Validation Report ⓘ

Nov 23, 2022 – 12:59 PM JST

PDB ID : 7V6C  
EMDB ID : EMD-31742  
Title : Structure of the Dicer-2-R2D2 heterodimer bound to small RNA duplex  
Authors : Yamaguchi, S.; Nishizawa, T.; Kusakizako, T.; Yamashita, K.; Tomita, A.; Hirano, H.; Nishimasu, H.; Nureki, O.  
Deposited on : 2021-08-20  
Resolution : 3.30 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:

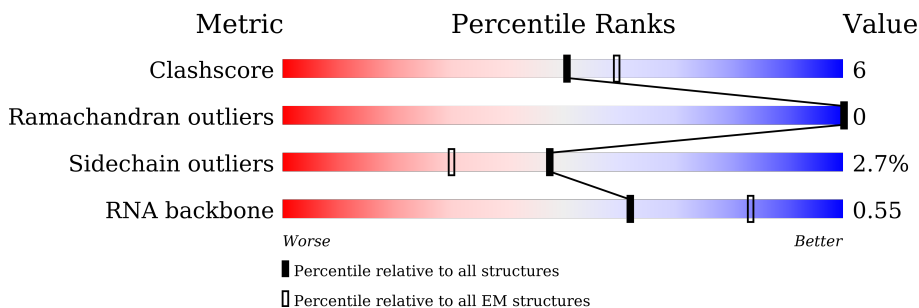
EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.3

# 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 3.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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	1733	
2	B	352	
3	E	21	
4	F	21	
5	C	22	
6	D	22	

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 15643 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 Dicer-2, isoform A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1540	12487	8033	2114	2273	67	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	HIS	-	expression tag	UNP A1ZAW0
A	-9	HIS	-	expression tag	UNP A1ZAW0
A	-8	HIS	-	expression tag	UNP A1ZAW0
A	-7	HIS	-	expression tag	UNP A1ZAW0
A	-6	HIS	-	expression tag	UNP A1ZAW0
A	-5	HIS	-	expression tag	UNP A1ZAW0
A	-4	HIS	-	expression tag	UNP A1ZAW0
A	-3	HIS	-	expression tag	UNP A1ZAW0
A	-2	ALA	-	expression tag	UNP A1ZAW0
A	-1	ALA	-	expression tag	UNP A1ZAW0
A	0	ALA	-	expression tag	UNP A1ZAW0
A	208	LEU	MET	engineered mutation	UNP A1ZAW0

- Molecule 2 is a protein called R2D2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	260	2057	1293	352	395	17	0	0

There are 41 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-40	MET	-	initiating methionine	UNP Q9VLW8
B	-39	ASP	-	expression tag	UNP Q9VLW8
B	-38	TYR	-	expression tag	UNP Q9VLW8
B	-37	LYS	-	expression tag	UNP Q9VLW8
B	-36	ASP	-	expression tag	UNP Q9VLW8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-35	HIS	-	expression tag	UNP Q9VLW8
B	-34	ASP	-	expression tag	UNP Q9VLW8
B	-33	GLY	-	expression tag	UNP Q9VLW8
B	-32	ASP	-	expression tag	UNP Q9VLW8
B	-31	TYR	-	expression tag	UNP Q9VLW8
B	-30	LYS	-	expression tag	UNP Q9VLW8
B	-29	ASP	-	expression tag	UNP Q9VLW8
B	-28	HIS	-	expression tag	UNP Q9VLW8
B	-27	ASP	-	expression tag	UNP Q9VLW8
B	-26	ILE	-	expression tag	UNP Q9VLW8
B	-25	ASP	-	expression tag	UNP Q9VLW8
B	-24	TYR	-	expression tag	UNP Q9VLW8
B	-23	LYS	-	expression tag	UNP Q9VLW8
B	-22	ASP	-	expression tag	UNP Q9VLW8
B	-21	ASP	-	expression tag	UNP Q9VLW8
B	-20	ASP	-	expression tag	UNP Q9VLW8
B	-19	ASP	-	expression tag	UNP Q9VLW8
B	-18	LYS	-	expression tag	UNP Q9VLW8
B	-17	HIS	-	expression tag	UNP Q9VLW8
B	-16	ARG	-	expression tag	UNP Q9VLW8
B	-15	TYR	-	expression tag	UNP Q9VLW8
B	-14	THR	-	expression tag	UNP Q9VLW8
B	-13	SER	-	expression tag	UNP Q9VLW8
B	-12	LEU	-	expression tag	UNP Q9VLW8
B	-11	TYR	-	expression tag	UNP Q9VLW8
B	-10	LYS	-	expression tag	UNP Q9VLW8
B	-9	LYS	-	expression tag	UNP Q9VLW8
B	-8	ALA	-	expression tag	UNP Q9VLW8
B	-7	GLY	-	expression tag	UNP Q9VLW8
B	-6	SER	-	expression tag	UNP Q9VLW8
B	-5	ALA	-	expression tag	UNP Q9VLW8
B	-4	ALA	-	expression tag	UNP Q9VLW8
B	-3	ALA	-	expression tag	UNP Q9VLW8
B	-2	PRO	-	expression tag	UNP Q9VLW8
B	-1	PHE	-	expression tag	UNP Q9VLW8
B	0	THR	-	expression tag	UNP Q9VLW8

- Molecule 3 is a RNA chain called RNA (5'-R(P\*UP\*GP\*AP\*GP\*G)-3').

Mol	Chain	Residues	Atoms				AltConf	Trace	
			Total	C	N	O			P
3	E	5	111	49	22	35	5	0	0

- Molecule 4 is a RNA chain called RNA (5'-R(P\*CP\*CP\*UP\*CP\*UP\*CP\*U)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
4	F	7	140	63	18	52	7	0	0

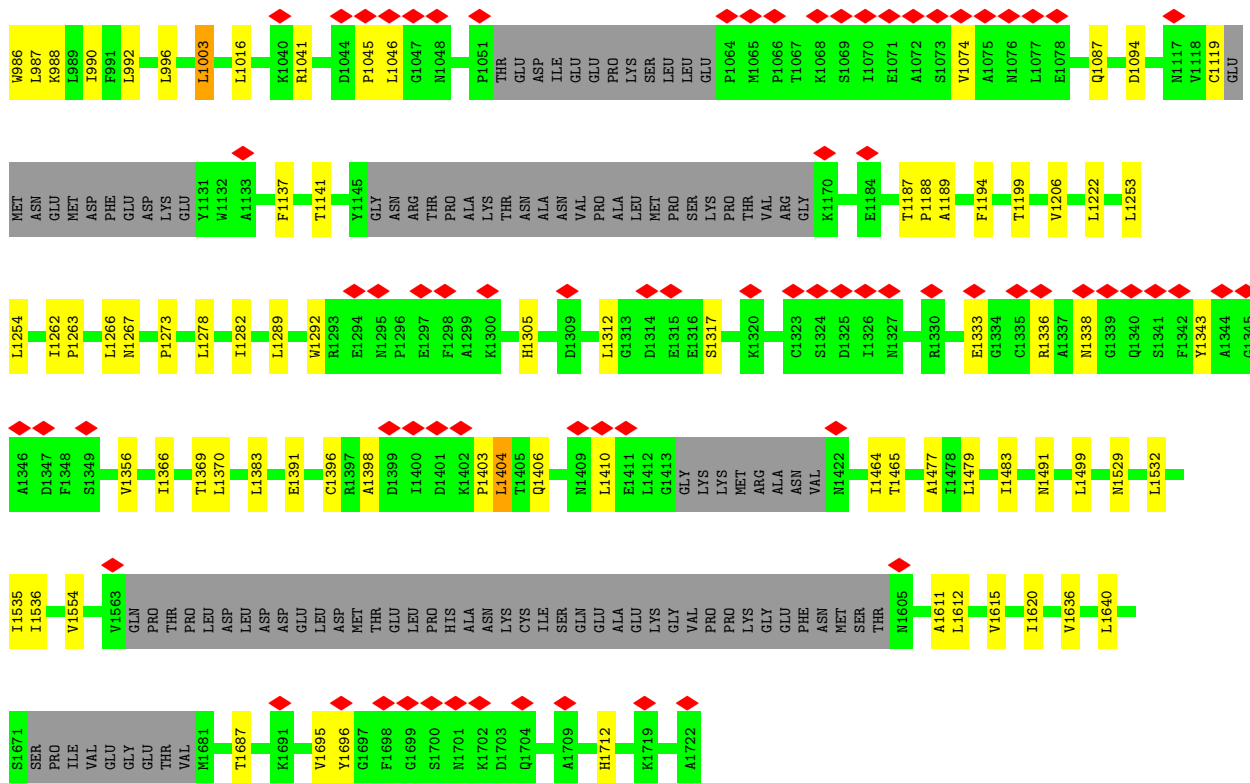
- Molecule 5 is a RNA chain called RNA (5'-R(\*AP\*UP\*GP\*AP\*GP\*GP\*UP\*AP\*GP\*UP\*AP\*GP\*GP\*UP\*UP\*GP\*UP\*AP\*UP\*AP\*G)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
5	C	21	435	193	79	143	20	0	1

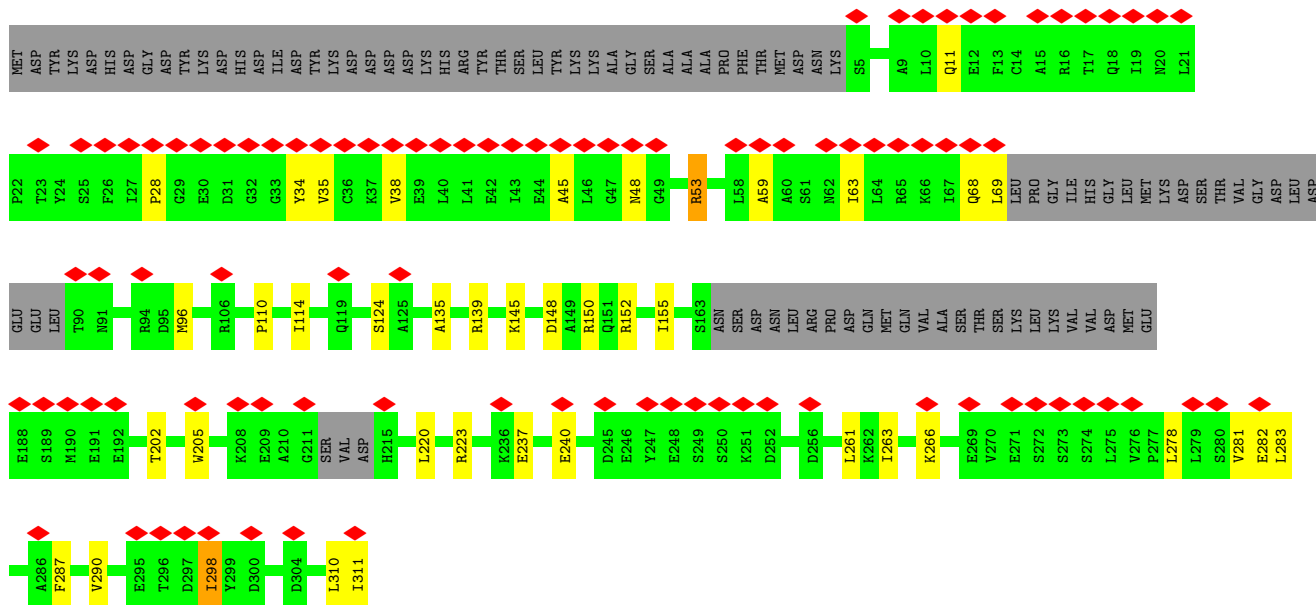
- Molecule 6 is a RNA chain called RNA (5'-R(\*CP\*UP\*AP\*UP\*AP\*CP\*AP\*AP\*CP\*CP\*UP\*AP\*CP\*UP\*AP\*CP\*CP\*UP\*CP\*UP\*C)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
6	D	21	413	186	66	141	20	0	1





• Molecule 2: R2D2



• Molecule 3: RNA (5'-R(P\*UP\*GP\*AP\*GP\*G)-3')

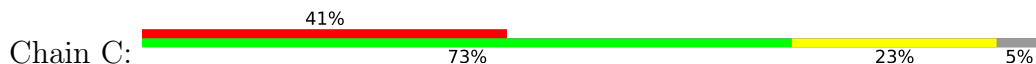




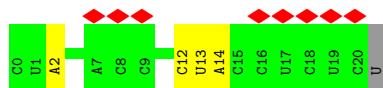
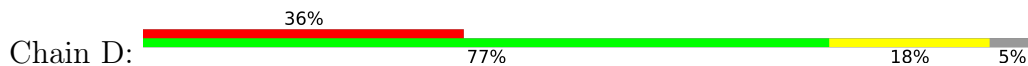
- Molecule 4: RNA (5'-R(P\*CP\*CP\*UP\*CP\*UP\*CP\*U)-3')



- Molecule 5: RNA (5'-R(\*AP\*UP\*GP\*AP\*GP\*GP\*UP\*AP\*GP\*UP\*AP\*GP\*GP\*UP\*UP\*GP\*UP\*AP\*UP\*AP\*G)-3')



- Molecule 6: RNA (5'-R(\*CP\*UP\*AP\*UP\*AP\*CP\*AP\*AP\*CP\*CP\*UP\*AP\*CP\*UP\*AP\*CP\*CP\*UP\*CP\*UP\*C)-3')





## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	144979	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 ( $e^-/\text{\AA}^2$ )	48	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.121	Depositor
Minimum map value	-0.070	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.0208	Depositor
Map size (Å)	211.68117, 211.68117, 211.68117	wwPDB
Map dimensions	212, 212, 212	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.99849606, 0.99849606, 0.99849606	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	0.61	0/12760	0.70	0/17269
2	B	0.63	0/2086	0.70	0/2802
3	E	0.70	0/124	0.84	0/192
4	F	0.23	0/153	0.63	0/234
5	C	0.57	0/487	0.80	0/760
6	D	0.28	0/458	0.68	0/709
All	All	0.60	0/16068	0.70	0/21966

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	12487	0	12531	153	0
2	B	2057	0	2072	25	0
3	E	111	0	55	2	0
4	F	140	0	75	2	0
5	C	435	0	214	3	0
6	D	413	0	215	3	0
All	All	15643	0	15162	183	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 6.

All (183) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LEU:HD22	1:A:133:LEU:HD21	1.57	0.87
1:A:184:VAL:HG11	1:A:528:ILE:HG12	1.57	0.86
1:A:717:LEU:HD12	1:A:718:PRO:HD2	1.60	0.82
1:A:814:TRP:HH2	1:A:820:LEU:HD12	1.45	0.81
1:A:1479:LEU:HD12	1:A:1620:ILE:HD13	1.62	0.81
1:A:886:TYR:HE1	1:A:975:LEU:HB3	1.54	0.71
1:A:94:VAL:HG11	1:A:118:VAL:HG13	1.73	0.70
1:A:568:PRO:HB2	1:A:643:ILE:HG22	1.74	0.69
1:A:328:THR:HG21	2:B:290:VAL:H	1.58	0.68
1:A:504:LEU:HD12	1:A:504:LEU:O	1.94	0.67
1:A:61:LEU:HD22	1:A:133:LEU:CD2	2.24	0.67
1:A:1141:THR:HG22	1:A:1141:THR:O	1.93	0.67
2:B:38:VAL:HG23	2:B:59:ALA:HB1	1.76	0.66
1:A:606:PHE:HB2	1:A:610:ALA:HB2	1.79	0.65
2:B:263:ILE:HG21	2:B:283:LEU:HD23	1.79	0.65
1:A:949:LEU:HB3	1:A:1273:PRO:HG2	1.79	0.65
1:A:1222:LEU:HA	1:A:1477:ALA:HB1	1.80	0.63
2:B:150:ARG:NH1	6:D:2:A:OP1	2.31	0.62
1:A:925:TYR:HB3	1:A:928:ARG:HB2	1.81	0.62
1:A:1356:VAL:O	1:A:1356:VAL:HG12	1.99	0.62
1:A:945:LEU:HD22	1:A:987:LEU:HD21	1.82	0.62
1:A:1396:CYS:HB2	1:A:1404:LEU:HD11	1.82	0.62
1:A:497:THR:HG22	1:A:499:GLU:H	1.65	0.61
1:A:895:VAL:HG13	1:A:940:ILE:HG23	1.81	0.61
1:A:1333:GLU:HA	1:A:1336:ARG:HE	1.65	0.60
1:A:1695:VAL:HG11	1:A:1712:HIS:HB3	1.83	0.60
2:B:35:VAL:HB	2:B:48:ASN:HB3	1.85	0.59
1:A:905:LEU:HA	1:A:916:THR:HB	1.82	0.59
1:A:41:VAL:HG11	1:A:169:VAL:HG21	1.85	0.59
1:A:81:THR:HG23	1:A:83:PHE:H	1.68	0.59
1:A:412:THR:O	1:A:414:GLU:N	2.35	0.59
2:B:96:MET:HG3	2:B:155:ILE:HG12	1.84	0.59
1:A:67:VAL:HG13	1:A:89:VAL:HG11	1.84	0.58
1:A:557:PRO:HB3	1:A:567:LEU:HD23	1.86	0.58
1:A:771:PHE:HE1	1:A:1206:VAL:HG11	1.69	0.57
1:A:1403:PRO:HG2	1:A:1406:GLN:HB2	1.87	0.57
1:A:181:ILE:HD13	1:A:521:TYR:HD1	1.70	0.56
1:A:881:ILE:HD11	1:A:892:PRO:HB2	1.88	0.56
2:B:310:LEU:O	2:B:311:ILE:C	2.44	0.56
1:A:757:LEU:HD13	1:A:844:PHE:CZ	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:587:PHE:HE1	1:A:1535:ILE:HD12	1.69	0.56
1:A:757:LEU:HD13	1:A:844:PHE:HZ	1.71	0.56
1:A:253:ASN:O	1:A:253:ASN:OD1	2.23	0.55
1:A:802:PHE:CD2	1:A:849:MET:HA	2.41	0.55
1:A:339:GLN:HE22	2:B:311:ILE:HG21	1.72	0.55
2:B:53:ARG:HD2	2:B:145:LYS:HD2	1.89	0.55
1:A:1262:ILE:N	1:A:1263:PRO:HD2	2.21	0.55
1:A:53:ILE:HD11	1:A:132:SER:HB3	1.89	0.54
1:A:1636:VAL:O	1:A:1640:LEU:HD13	2.07	0.54
1:A:729:ILE:HD11	1:A:756:ILE:HG12	1.88	0.54
1:A:1391:GLU:HB2	1:A:1404:LEU:HD12	1.90	0.54
1:A:87:PHE:CZ	1:A:89:VAL:HG21	2.43	0.53
1:A:1536:ILE:HG12	1:A:1612:LEU:HD11	1.91	0.53
1:A:272:SER:HA	1:A:323:HIS:HE1	1.74	0.53
1:A:154:MET:O	1:A:157:PHE:HB2	2.08	0.53
1:A:808:ARG:HA	1:A:814:TRP:HB2	1.90	0.53
1:A:66:THR:HG22	1:A:68:GLU:H	1.74	0.53
1:A:934:HIS:HB3	1:A:937:LYS:HD3	1.91	0.52
1:A:886:TYR:CE1	1:A:975:LEU:HB3	2.41	0.52
2:B:266:LYS:HB3	2:B:282:GLU:HB3	1.90	0.52
1:A:881:ILE:HG23	1:A:979:PHE:HB2	1.92	0.52
4:F:15:C:H2'	4:F:16:C:C6	2.45	0.52
1:A:477:PHE:HB3	1:A:505:PHE:HE2	1.76	0.51
1:A:814:TRP:HH2	1:A:820:LEU:CD1	2.18	0.51
1:A:19:LEU:HD21	1:A:26:VAL:HG22	1.92	0.51
4:F:16:C:H2'	4:F:17:U:C6	2.46	0.51
2:B:148:ASP:O	2:B:152:ARG:HG2	2.12	0.50
1:A:881:ILE:HG21	1:A:986:TRP:CE3	2.47	0.50
1:A:900:ARG:HA	1:A:938:PHE:CG	2.46	0.50
1:A:802:PHE:HD2	1:A:849:MET:HA	1.77	0.50
1:A:1312:LEU:C	1:A:1312:LEU:HD12	2.31	0.49
1:A:286:TYR:HA	1:A:289:LYS:HE2	1.93	0.49
1:A:754:TYR:HE1	1:A:988:LYS:HB3	1.78	0.49
1:A:67:VAL:HG22	1:A:89:VAL:CG1	2.42	0.49
1:A:465:GLU:HG3	1:A:466:GLU:HG3	1.94	0.49
1:A:344:MET:HB2	1:A:362:PHE:HE2	1.77	0.49
1:A:124:THR:C	1:A:126:THR:H	2.15	0.49
1:A:754:TYR:HE2	1:A:992:LEU:HD13	1.78	0.49
1:A:222:MET:SD	1:A:504:LEU:HD11	2.53	0.48
1:A:608:VAL:HA	1:A:611:LYS:HE2	1.96	0.48
1:A:1398:ALA:HB2	1:A:1404:LEU:HG	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:ILE:HD11	1:A:333:ILE:HD13	1.94	0.48
1:A:61:LEU:HD12	1:A:112:LEU:HB2	1.95	0.48
2:B:261:LEU:HD11	2:B:310:LEU:HD11	1.95	0.48
1:A:727:TYR:CE2	1:A:785:PRO:HG3	2.49	0.48
1:A:731:LEU:HG	1:A:779:VAL:HG22	1.96	0.48
1:A:1187:THR:HG23	1:A:1188:PRO:HD2	1.96	0.48
1:A:59:ARG:HD2	1:A:110:GLN:HA	1.96	0.48
1:A:388:CYS:HB3	1:A:475:HIS:HB2	1.97	0.47
1:A:750:THR:HG21	1:A:984:ASP:OD2	2.14	0.47
1:A:860:GLY:HA2	1:A:885:TRP:CE3	2.49	0.47
2:B:110:PRO:CG	2:B:135:ALA:HB2	2.45	0.47
1:A:716:ALA:HB1	1:A:758:LEU:HB2	1.97	0.47
2:B:28:PRO:HA	2:B:34:TYR:HD1	1.80	0.47
1:A:58:LYS:HB3	1:A:134:SER:HB2	1.96	0.47
1:A:272:SER:HA	1:A:323:HIS:CE1	2.50	0.47
1:A:756:ILE:HD11	1:A:996:LEU:HD21	1.96	0.47
6:D:13:U:H2'	6:D:14:A:C8	2.50	0.47
1:A:1312:LEU:HD12	1:A:1317:SER:HB2	1.97	0.46
2:B:220:LEU:HD12	2:B:223:ARG:HD3	1.97	0.46
2:B:281:VAL:HG23	2:B:298:ILE:HD12	1.96	0.46
1:A:463:VAL:HG11	1:A:490:GLN:HG3	1.97	0.46
1:A:939:MET:HE2	1:A:973:PRO:HD3	1.98	0.46
1:A:1199:THR:HA	1:A:1267:ASN:HB3	1.97	0.46
1:A:1404:LEU:HD23	1:A:1404:LEU:HA	1.76	0.46
1:A:832:VAL:HG11	1:A:844:PHE:CD2	2.51	0.46
1:A:632:SER:HB3	1:A:645:ALA:HB2	1.98	0.45
2:B:124:SER:HB2	6:D:12:C:O2'	2.17	0.45
1:A:58:LYS:HA	1:A:132:SER:O	2.16	0.45
1:A:866:ARG:HH12	1:A:975:LEU:HD23	1.80	0.45
1:A:754:TYR:CE2	1:A:992:LEU:HD13	2.52	0.45
1:A:484:THR:HG22	1:A:486:ASN:H	1.81	0.45
1:A:1305:HIS:CE1	1:A:1343:TYR:CD2	3.05	0.45
1:A:1338:ASN:OD1	1:A:1343:TYR:HB2	2.16	0.45
1:A:12:GLN:O	1:A:16:VAL:HG23	2.16	0.45
1:A:33:GLY:O	1:A:37:VAL:HG23	2.17	0.45
1:A:1312:LEU:CD1	1:A:1317:SER:HB2	2.46	0.45
1:A:239:SER:O	1:A:243:GLU:HG2	2.15	0.45
1:A:242:ILE:HG22	1:A:246:TYR:CE2	2.52	0.45
1:A:506:THR:HG22	1:A:507:ALA:N	2.31	0.45
1:A:199:ILE:O	1:A:535:ARG:HB3	2.17	0.45
1:A:1289:LEU:HD12	1:A:1289:LEU:HA	1.87	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1499:LEU:HD23	1:A:1499:LEU:HA	1.79	0.44
1:A:986:TRP:O	1:A:990:ILE:HG13	2.17	0.44
1:A:1410:LEU:HD12	1:A:1410:LEU:HA	1.79	0.44
1:A:38:ALA:HB2	1:A:171:LEU:HD11	2.00	0.44
1:A:953:VAL:O	1:A:953:VAL:HG22	2.17	0.44
2:B:69:LEU:C	2:B:69:LEU:HD12	2.38	0.44
1:A:15:LEU:HD22	1:A:26:VAL:HG13	1.99	0.44
1:A:116:ALA:HB2	1:A:150:PHE:CE1	2.53	0.44
1:A:124:THR:C	1:A:126:THR:N	2.71	0.44
1:A:276:PHE:CE2	1:A:310:LYS:HG3	2.53	0.44
1:A:89:VAL:HG12	1:A:90:GLY:N	2.33	0.44
1:A:1045:PRO:O	1:A:1046:LEU:HB2	2.18	0.44
1:A:19:LEU:HD13	1:A:19:LEU:HA	1.86	0.43
1:A:283:ASP:O	1:A:287:GLN:HG2	2.18	0.43
1:A:808:ARG:HG3	1:A:814:TRP:CG	2.53	0.43
2:B:110:PRO:HG3	2:B:135:ALA:HB2	1.98	0.43
2:B:38:VAL:HG11	2:B:63:ILE:HG21	2.00	0.43
5:C:14:U:H2'	5:C:15:G:C8	2.54	0.43
1:A:568:PRO:O	1:A:571:ALA:HB3	2.18	0.43
1:A:932:VAL:HG23	1:A:935:LYS:HD2	2.01	0.43
1:A:378:PHE:CD2	1:A:419:LEU:HD11	2.53	0.43
1:A:754:TYR:CE1	1:A:988:LYS:HB3	2.54	0.43
2:B:68:GLN:C	2:B:69:LEU:HG	2.39	0.43
1:A:1253:LEU:HD22	1:A:1370:LEU:HD22	2.01	0.43
1:A:388:CYS:SG	1:A:457:LEU:HD12	2.59	0.42
2:B:237:GLU:HA	2:B:240:GLU:HG2	2.00	0.42
1:A:388:CYS:HB2	1:A:475:HIS:O	2.19	0.42
1:A:504:LEU:HD12	1:A:504:LEU:C	2.39	0.42
1:A:115:THR:HG22	1:A:116:ALA:N	2.34	0.42
2:B:45:ALA:HB2	2:B:63:ILE:HD13	2.00	0.42
1:A:1278:LEU:HD23	1:A:1278:LEU:H	1.84	0.42
3:E:1:U:H2'	3:E:2:G:C8	2.55	0.42
1:A:436:SER:O	1:A:437:VAL:C	2.58	0.42
1:A:1312:LEU:HD12	1:A:1312:LEU:O	2.20	0.42
1:A:175:LEU:HD23	1:A:175:LEU:HA	1.95	0.41
1:A:717:LEU:HB2	1:A:828:SER:HB3	2.01	0.41
1:A:518:ILE:O	1:A:522:ARG:HG3	2.21	0.41
3:E:2:G:H2'	3:E:3:A:C8	2.55	0.41
1:A:1189:ALA:HB1	1:A:1194:PHE:HE2	1.85	0.41
1:A:144:GLY:O	1:A:151:ARG:HB2	2.21	0.41
1:A:1266:LEU:HD21	1:A:1369:THR:HG21	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1611:ALA:O	1:A:1615:VAL:HG23	2.21	0.41
5:C:1:U:H2'	5:C:2:G:C8	2.54	0.41
1:A:252:MET:O	1:A:274:LYS:HE3	2.20	0.41
1:A:392:VAL:O	1:A:462:SER:N	2.53	0.41
1:A:401:ILE:HD13	1:A:401:ILE:HA	1.97	0.41
1:A:1041:ARG:HD3	1:A:1041:ARG:HA	1.77	0.41
1:A:1464:ILE:HG22	1:A:1465:THR:HG23	2.03	0.41
1:A:67:VAL:HG22	1:A:89:VAL:HG13	2.03	0.41
1:A:1479:LEU:HD21	1:A:1483:ILE:HD12	2.02	0.41
1:A:1529:ASN:HB3	1:A:1532:LEU:HB3	2.03	0.41
1:A:19:LEU:HD12	1:A:169:VAL:HG22	2.03	0.40
1:A:562:ASN:HD21	1:A:661:GLU:HB3	1.87	0.40
2:B:145:LYS:HG3	5:C:11:G:OP1	2.21	0.40
1:A:765:LEU:HD21	1:A:1003:LEU:HD12	2.04	0.40
1:A:813:ILE:HG23	1:A:1273:PRO:O	2.20	0.40
1:A:1254:LEU:HD12	1:A:1366:ILE:HG21	2.03	0.40
1:A:1491:ASN:HB2	1:A:1499:LEU:HD21	2.03	0.40
2:B:139:ARG:HD3	2:B:139:ARG:HA	1.78	0.40
1:A:324:ARG:O	1:A:328:THR:HG23	2.22	0.40
1:A:405:LEU:HD23	1:A:405:LEU:HA	1.91	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	A	1514/1733 (87%)	1461 (96%)	53 (4%)	0	100	100
2	B	252/352 (72%)	244 (97%)	8 (3%)	0	100	100
All	All	1766/2085 (85%)	1705 (96%)	61 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	1385/1557 (89%)	1349 (97%)	36 (3%)	46 71
2	B	230/312 (74%)	222 (96%)	8 (4%)	36 64
All	All	1615/1869 (86%)	1571 (97%)	44 (3%)	48 71

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	61	LEU
1	A	81	THR
1	A	96	ASP
1	A	141	CYS
1	A	159	ILE
1	A	205	THR
1	A	210	ASN
1	A	216	THR
1	A	389	LEU
1	A	400	CYS
1	A	404	LEU
1	A	418	VAL
1	A	466	GLU
1	A	628	ASP
1	A	633	ASP
1	A	881	ILE
1	A	913	GLN
1	A	925	TYR
1	A	942	VAL
1	A	949	LEU
1	A	952	TYR
1	A	1003	LEU
1	A	1016	LEU
1	A	1074	VAL
1	A	1087	GLN
1	A	1094	ASP

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Mol	Chain	Res	Type
1	A	1119	CYS
1	A	1137	PHE
1	A	1282	ILE
1	A	1292	TRP
1	A	1383	LEU
1	A	1404	LEU
1	A	1554	VAL
1	A	1687	THR
1	A	1696	TYR
2	B	11	GLN
2	B	53	ARG
2	B	114	ILE
2	B	202	THR
2	B	205	TRP
2	B	278	LEU
2	B	287	PHE
2	B	298	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	ASN
1	A	323	HIS
1	A	339	GLN
1	A	354	ASN
1	A	407	ASN
1	A	520	GLN
1	A	562	ASN
1	A	619	ASN
1	A	623	ASN
1	A	660	ASN
1	A	734	GLN
1	A	841	GLN
1	A	859	HIS
1	A	884	GLN
1	A	888	ASN
1	A	1098	ASN
1	A	1287	ASN
2	B	11	GLN
2	B	91	ASN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	E	4/21 (19%)	0	0
4	F	6/21 (28%)	0	0
5	C	19/22 (86%)	0	0
6	D	19/22 (86%)	0	0
All	All	48/86 (55%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

### 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 monosaccharides 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.

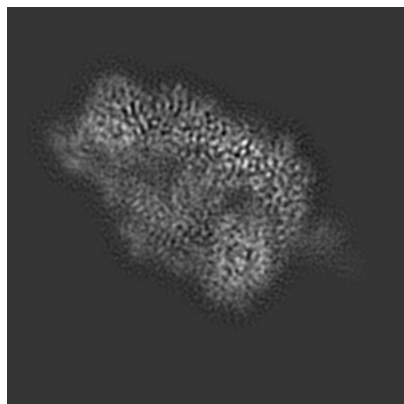
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31742. These allow visual inspection of the internal detail of the map and identification of artifacts.

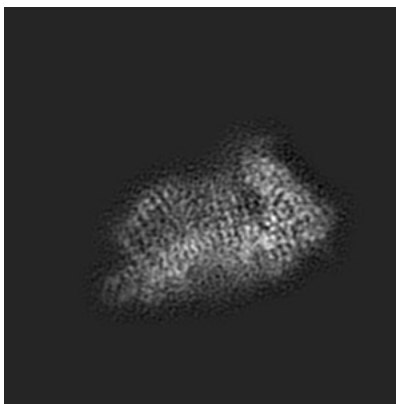
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

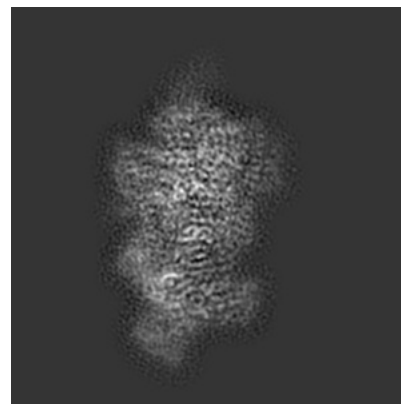
#### 6.1.1 Primary map



X

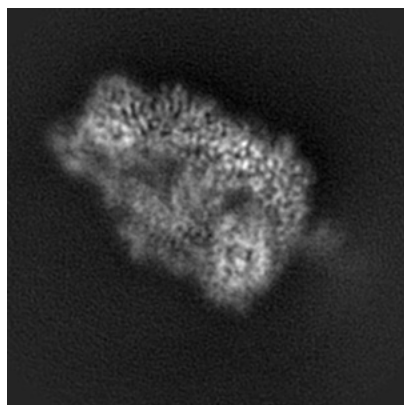


Y

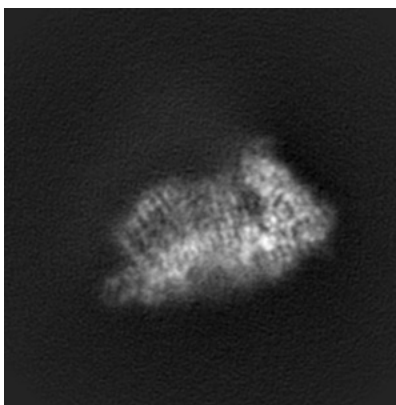


Z

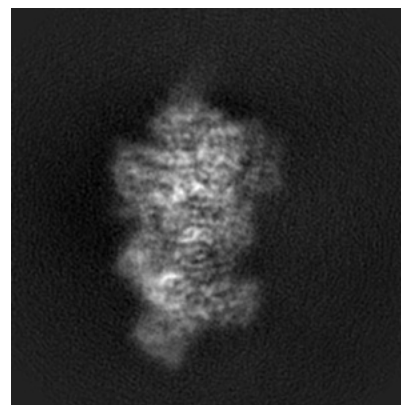
#### 6.1.2 Raw map



X



Y

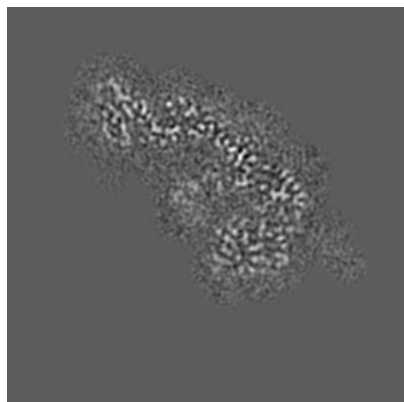


Z

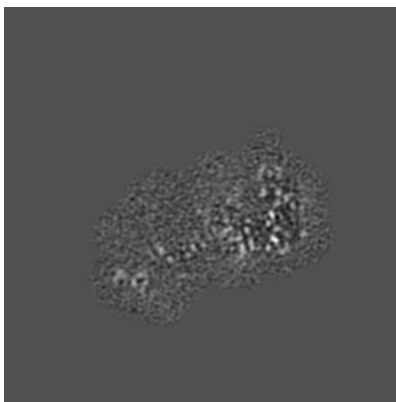
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

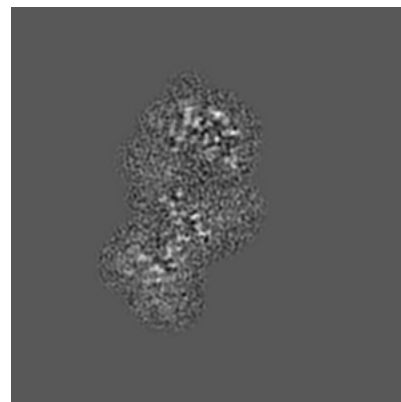
### 6.2.1 Primary map



X Index: 106

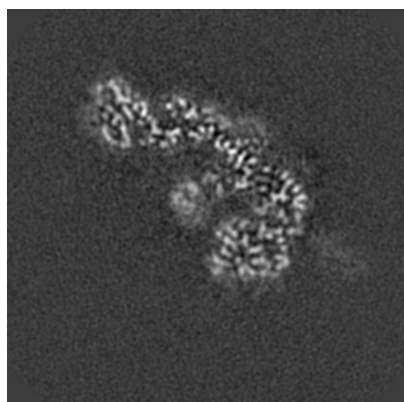


Y Index: 106

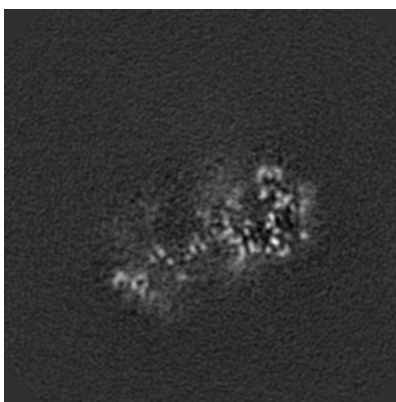


Z Index: 106

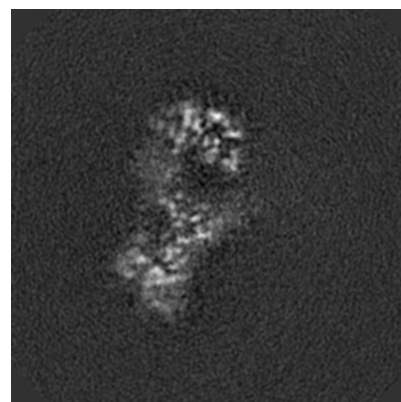
### 6.2.2 Raw map



X Index: 106



Y Index: 106

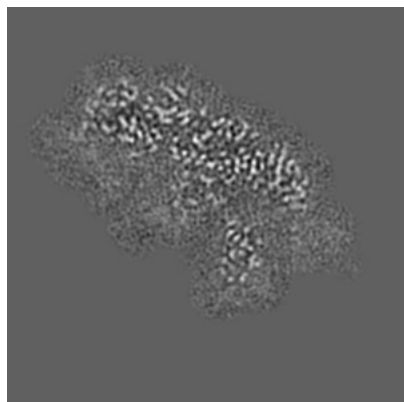


Z Index: 106

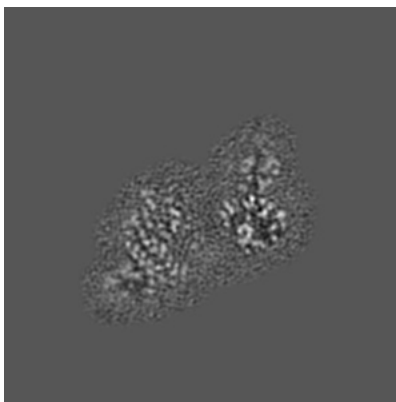
The images above show central slices of the map in three orthogonal directions.

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

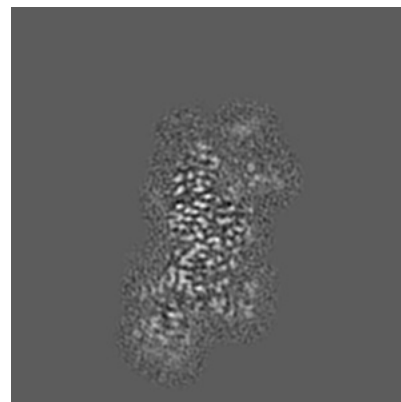
### 6.3.1 Primary map



X Index: 94

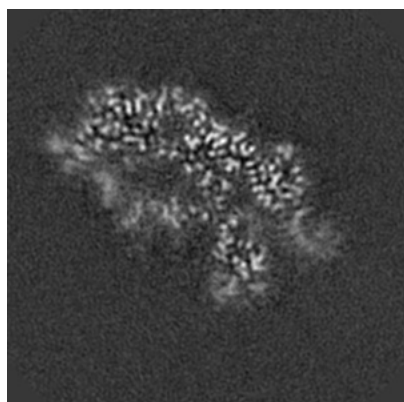


Y Index: 122

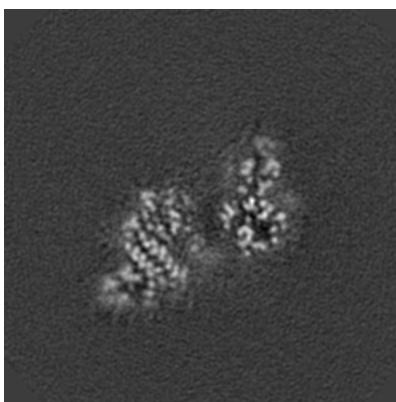


Z Index: 142

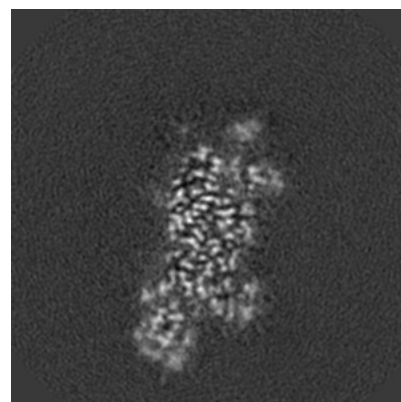
### 6.3.2 Raw map



X Index: 89



Y Index: 122

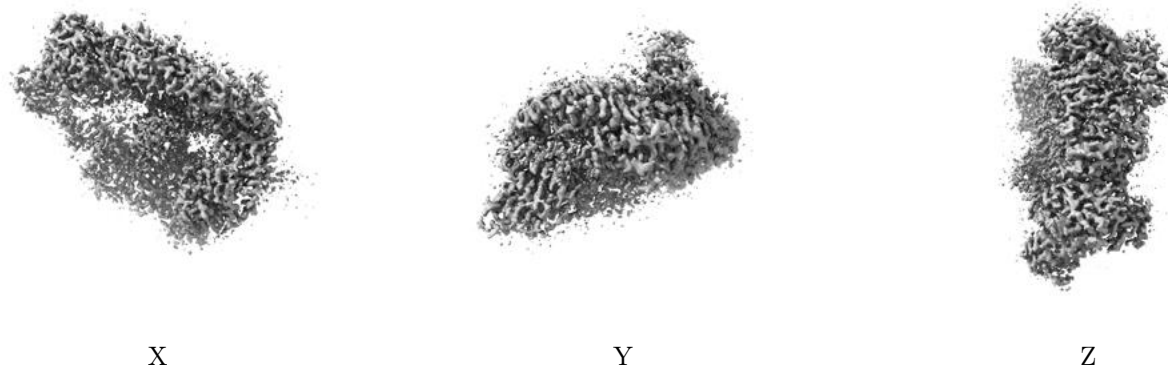


Z Index: 141

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

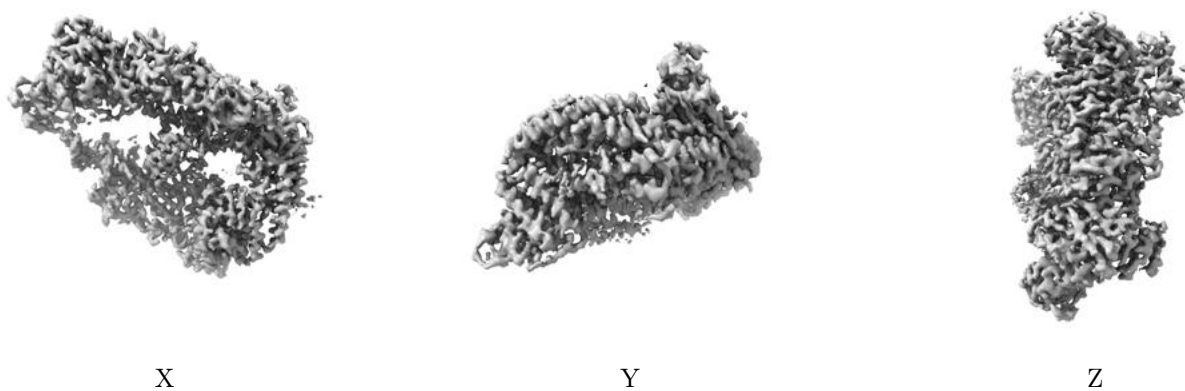
## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0208. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.4.2 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.

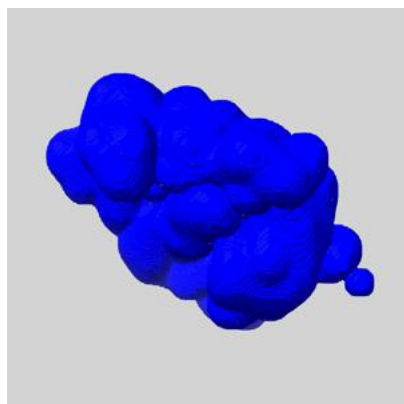
## 6.5 Mask visualisation [i](#)

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

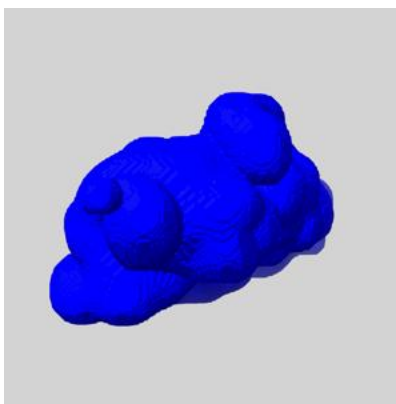
A mask typically either:

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

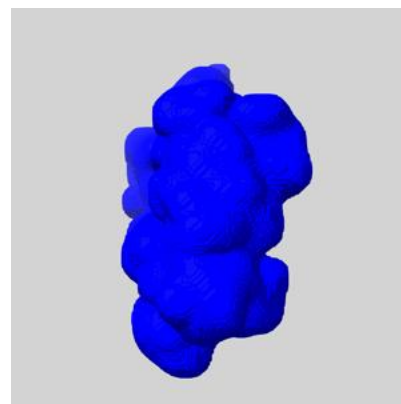
### 6.5.1 emd\_31742\_msk\_1.map [i](#)



X



Y

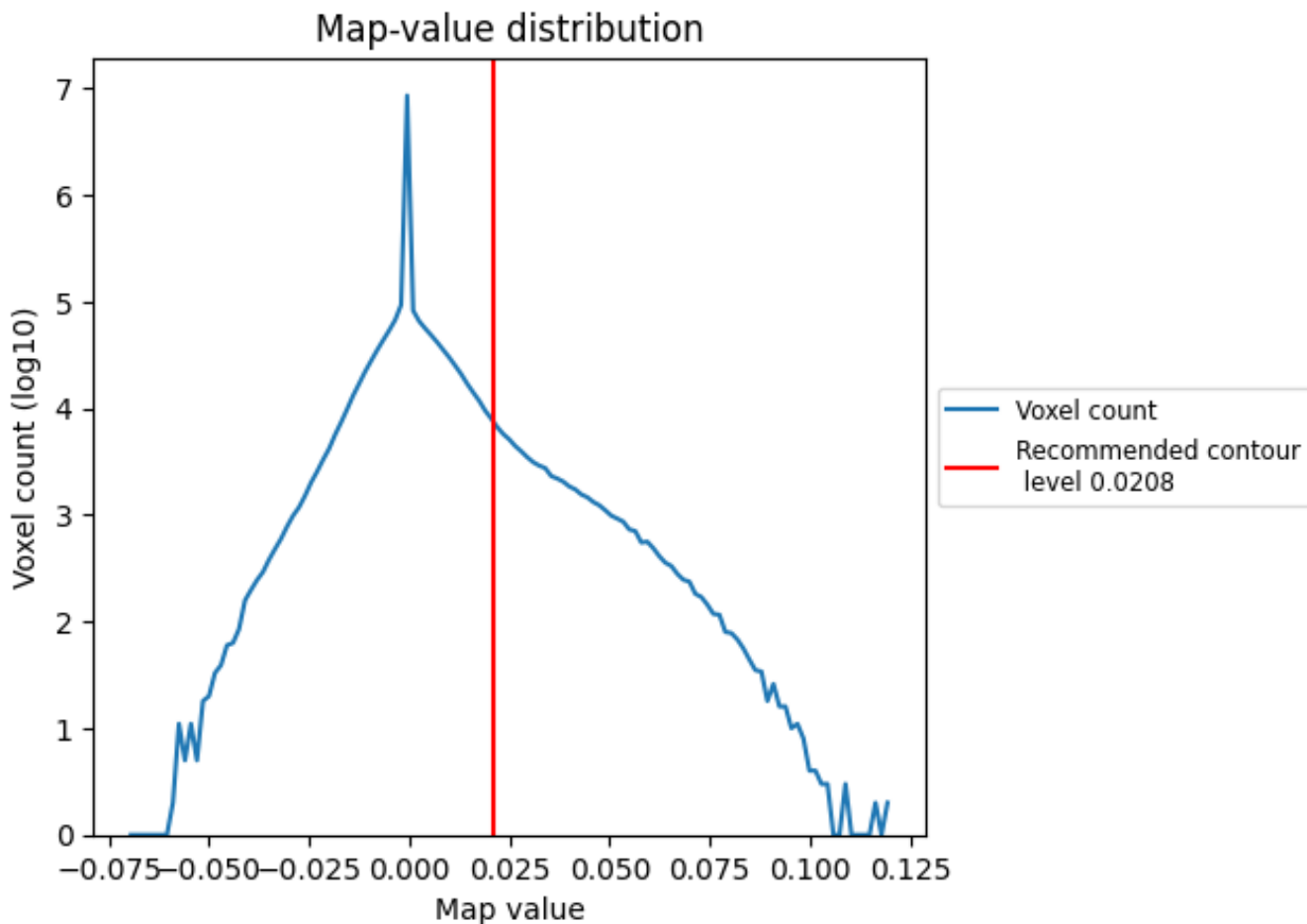


Z

## 7 Map analysis [i](#)

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

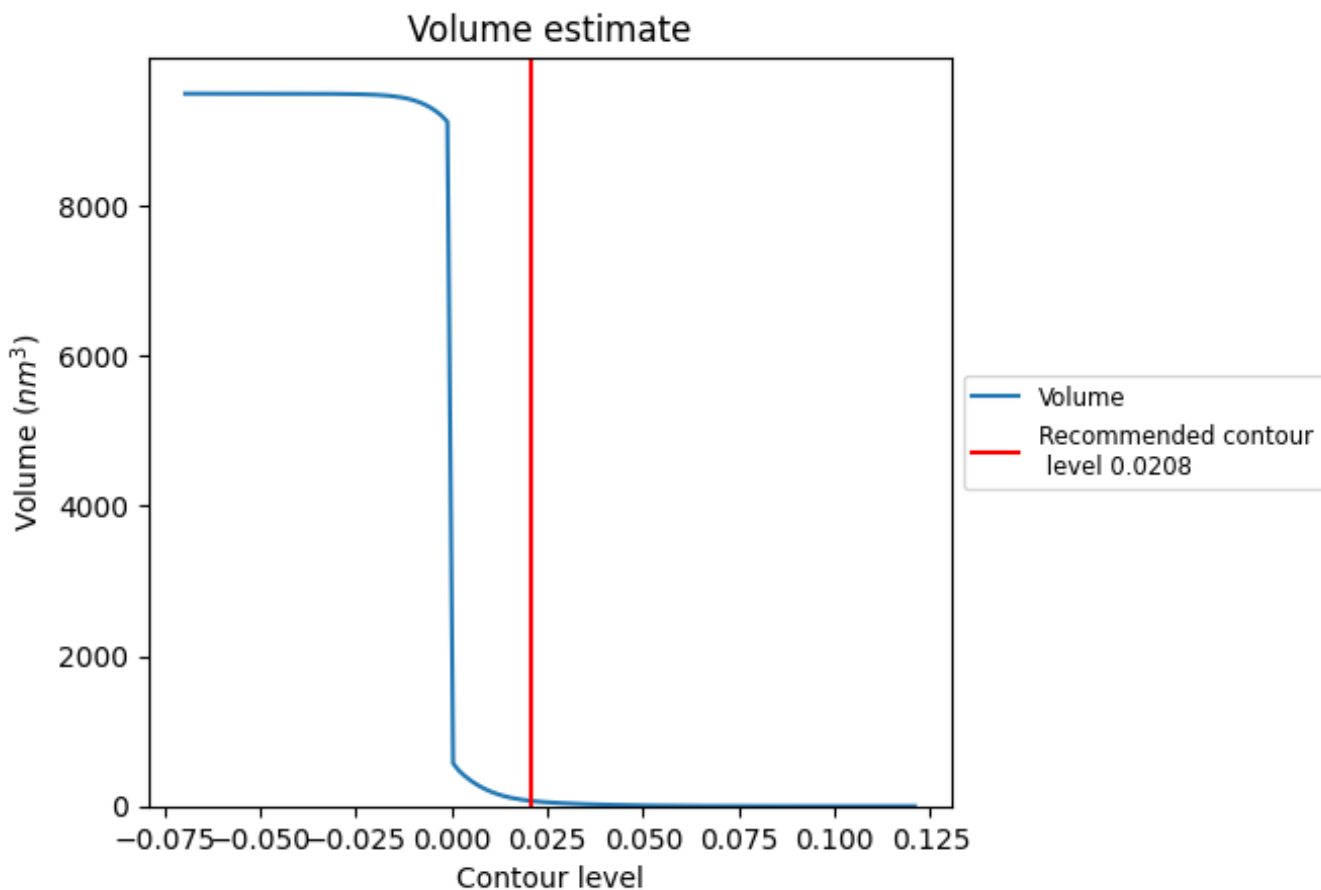
### 7.1 Map-value distribution [i](#)



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.



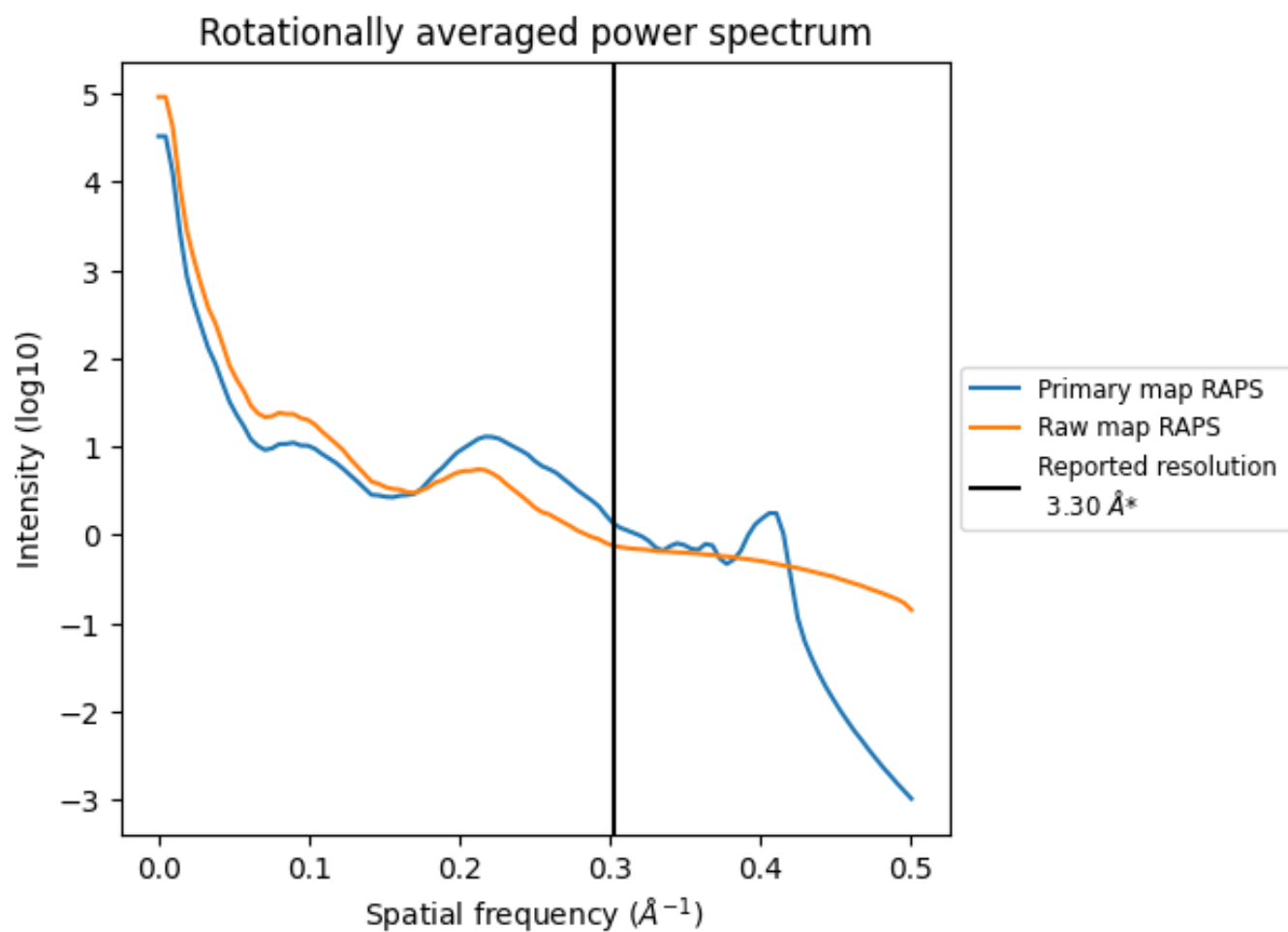
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is  $70 \text{ nm}^3$ ; this corresponds to an approximate mass of 63 kDa.

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

### 7.3 Rotationally averaged power spectrum i

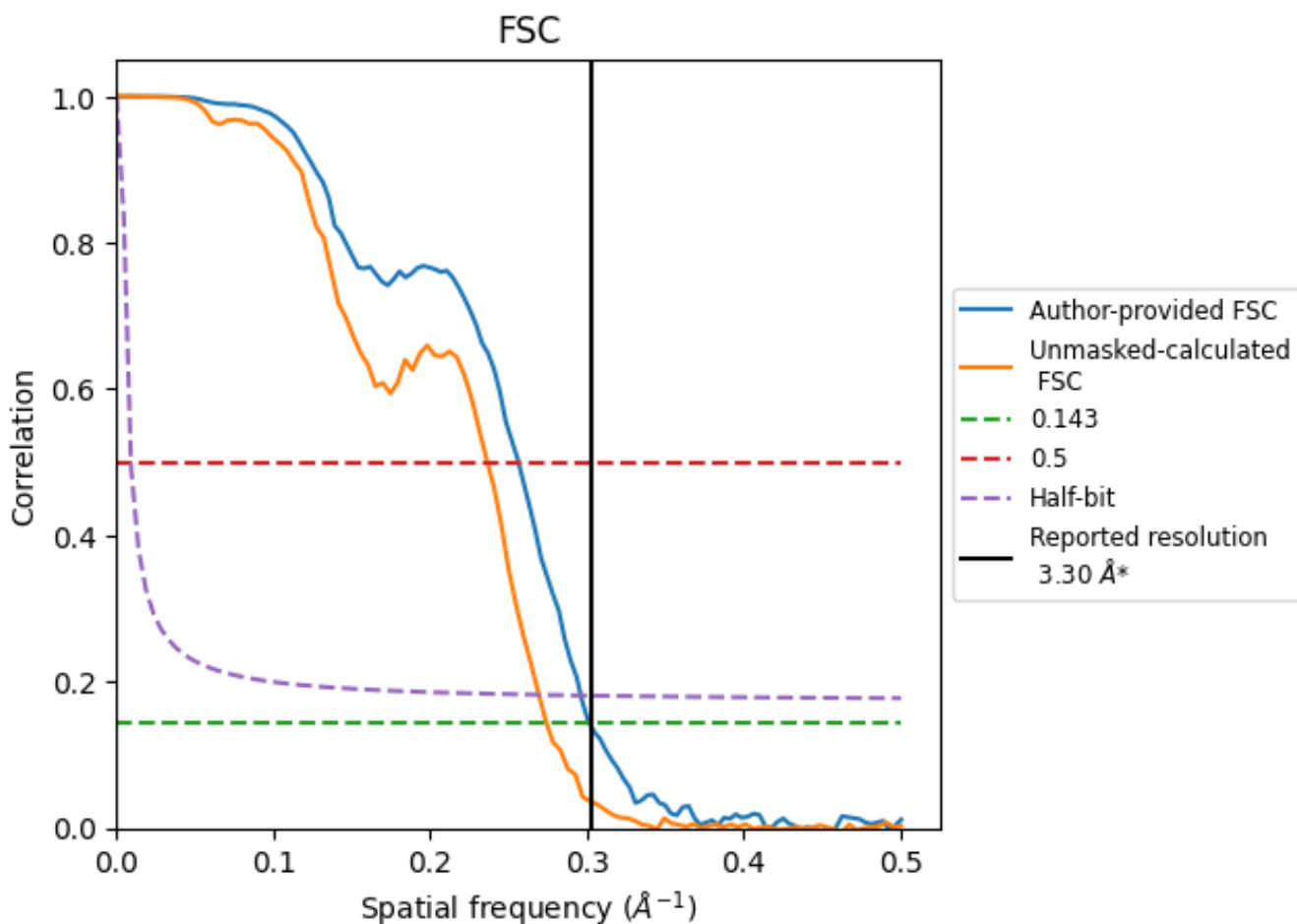


\*Reported resolution corresponds to spatial frequency of 0.303 Å<sup>-1</sup>

## 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 0.303 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

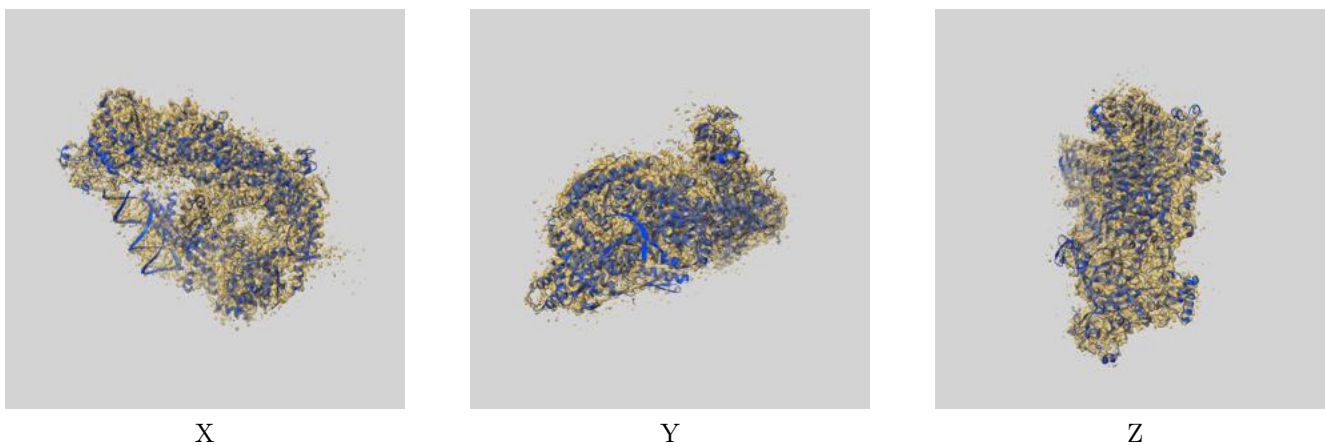
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	3.31	3.90	3.37
Unmasked-calculated*	3.64	4.22	3.71

\*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 3.64 differs from the reported value 3.3 by more than 10 %

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

This section contains information regarding the fit between EMDB map EMD-31742 and PDB model 7V6C. Per-residue inclusion information can be found in section 3 on page 6.

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



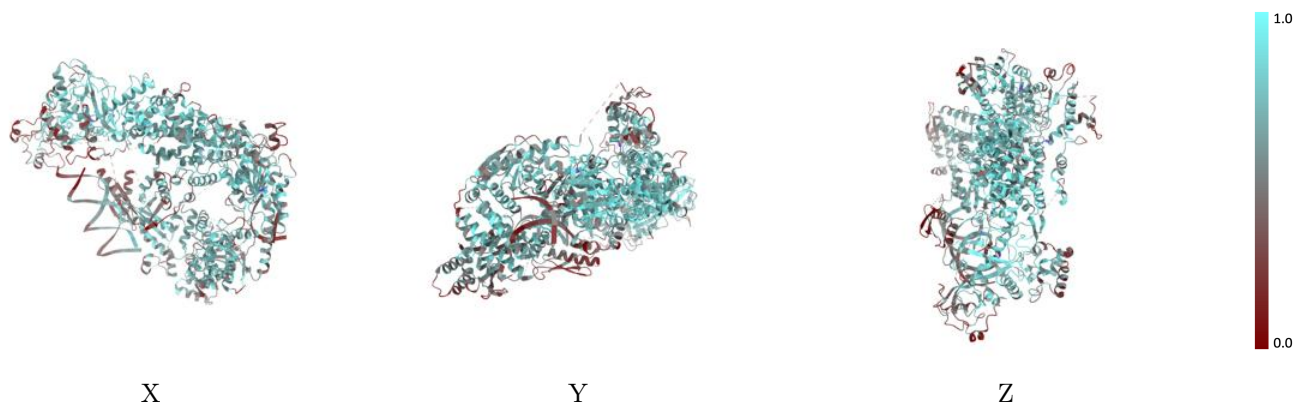
The images above show the 3D surface view of the map at the recommended contour level 0.0208 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



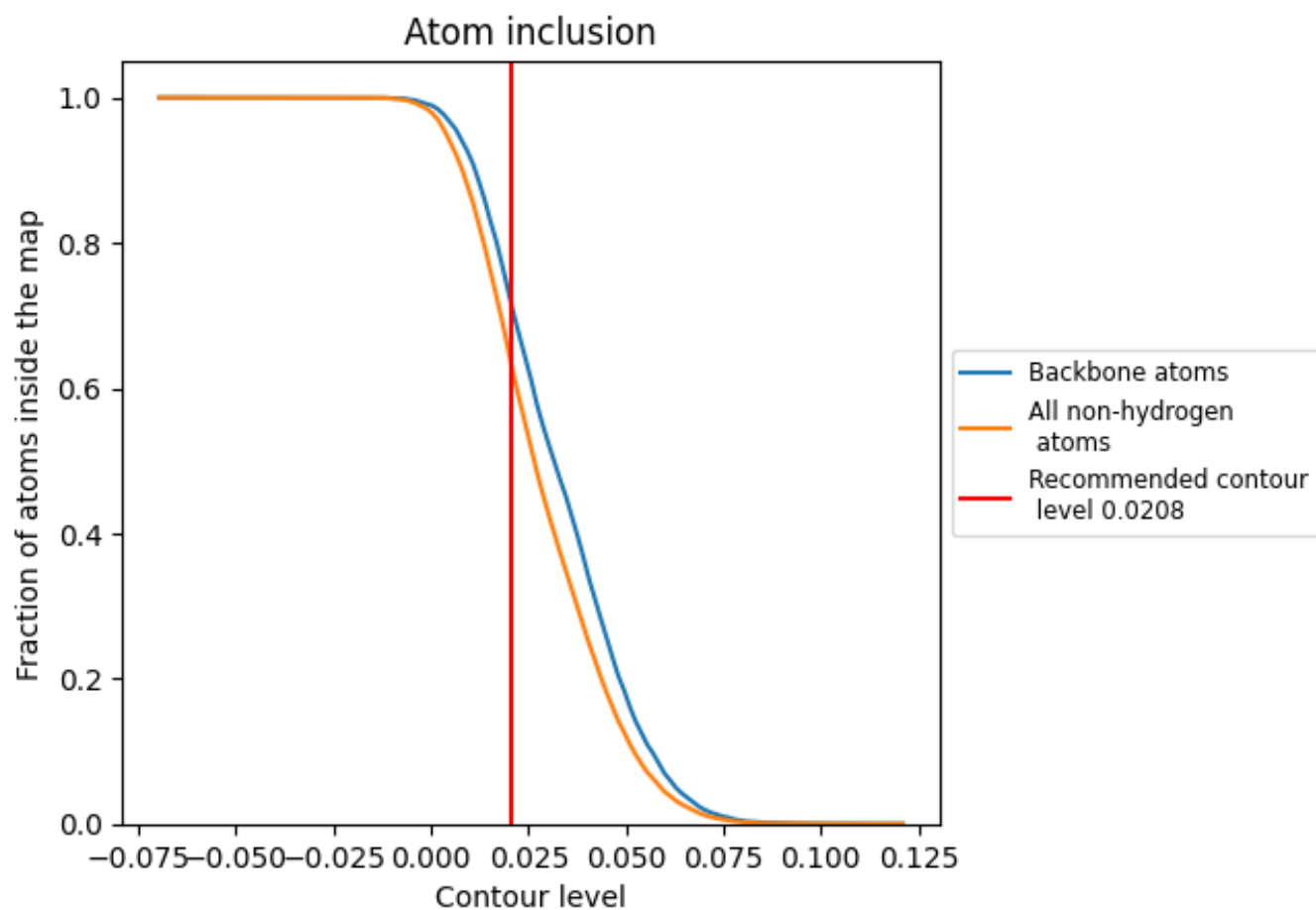
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0208).















## 9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 63% of all non-hydrogen atoms, are inside the map.

## 9.5 Map-model fit summary [i](#)

The table lists the average atom inclusion at the recommended contour level (0.0208) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6276	 0.4730
A	 0.6664	 0.4920
B	 0.5000	 0.4240
C	 0.4644	 0.3360
D	 0.5109	 0.3880
E	 0.2523	 0.2850
F	 0.2214	 0.2640

