



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 09:33 PM UTC

PDB ID : 9ECV / pdb_00009ecv
EMDB ID : EMD-47927
Title : CryoEM Structure Of Respiratory Syncytial Virus Polymerase in complex with Novel Non-Nucleoside Inhibitor Compound 16
Authors : Yin, Y.; Tran, M.T.; Yu, X.; Jonckers, T.; Carney, C.
Deposited on : 2024-11-15
Resolution : 2.79 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:

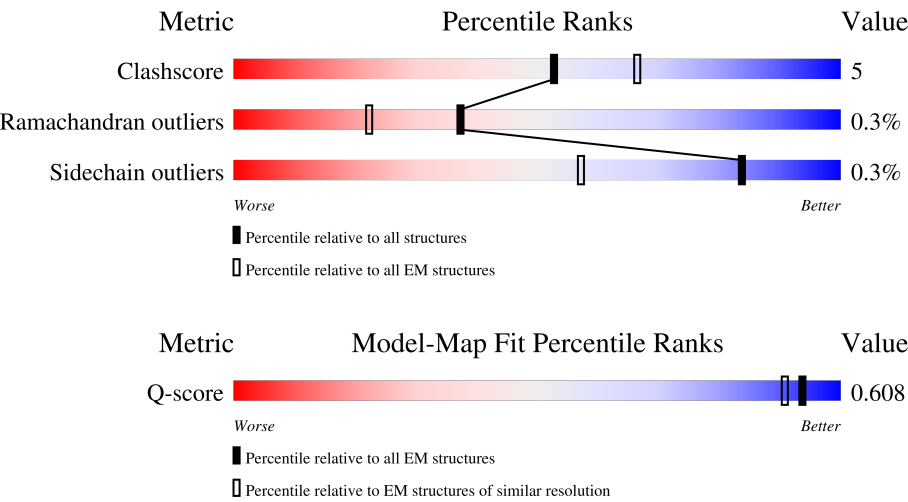
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.79 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	10811 (2.29 - 3.29)

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	2201	<div><div>53% 9% 38%</div><div>21%</div></div>
2	B	256	<div><div>37% 7% 55%</div><div>21%</div></div>
2	C	256	<div><div>18% 6% 76%</div><div>14%</div></div>
2	D	256	<div><div>19% 5% 76%</div><div>14%</div></div>

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Mol	Chain	Length	Quality of chain
2	E	256	<div><div><div></div><div></div><div></div></div><div>19%20%6%73%</div></div>
3	G	3	<div><div><div></div><div></div><div></div></div><div>33%33%33%</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13558 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 RNA-directed RNA polymerase L.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1367	Total	C	N	O	S	0	0
			11153	7186	1872	2038	57		

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-35	MET	-	initiating methionine	UNP P28887
A	-34	GLY	-	expression tag	UNP P28887
A	-33	SER	-	expression tag	UNP P28887
A	-32	TRP	-	expression tag	UNP P28887
A	-31	SER	-	expression tag	UNP P28887
A	-30	HIS	-	expression tag	UNP P28887
A	-29	PRO	-	expression tag	UNP P28887
A	-28	GLN	-	expression tag	UNP P28887
A	-27	PHE	-	expression tag	UNP P28887
A	-26	GLU	-	expression tag	UNP P28887
A	-25	LYS	-	expression tag	UNP P28887
A	-24	GLY	-	expression tag	UNP P28887
A	-23	SER	-	expression tag	UNP P28887
A	-22	GLY	-	expression tag	UNP P28887
A	-21	SER	-	expression tag	UNP P28887
A	-20	GLY	-	expression tag	UNP P28887
A	-19	SER	-	expression tag	UNP P28887
A	-18	SER	-	expression tag	UNP P28887
A	-17	TRP	-	expression tag	UNP P28887
A	-16	SER	-	expression tag	UNP P28887
A	-15	HIS	-	expression tag	UNP P28887
A	-14	PRO	-	expression tag	UNP P28887
A	-13	GLN	-	expression tag	UNP P28887
A	-12	PHE	-	expression tag	UNP P28887
A	-11	GLU	-	expression tag	UNP P28887
A	-10	LYS	-	expression tag	UNP P28887
A	-9	GLY	-	expression tag	UNP P28887
A	-8	SER	-	expression tag	UNP P28887

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	expression tag	UNP P28887
A	-6	SER	-	expression tag	UNP P28887
A	-5	LEU	-	expression tag	UNP P28887
A	-4	VAL	-	expression tag	UNP P28887
A	-3	PRO	-	expression tag	UNP P28887
A	-2	ARG	-	expression tag	UNP P28887
A	-1	GLY	-	expression tag	UNP P28887
A	0	SER	-	expression tag	UNP P28887

- Molecule 2 is a protein called Phosphoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	114	Total	C	N	O	S	0	0
			885	531	157	191	6		
2	C	62	Total	C	N	O	S	0	0
			478	291	89	94	4		
2	D	61	Total	C	N	O	S	0	0
			467	285	85	93	4		
2	E	69	Total	C	N	O	S	0	0
			535	327	101	102	5		

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	171	VAL	ILE	variant	UNP P03421
B	242	LYS	-	expression tag	UNP P03421
B	243	GLY	-	expression tag	UNP P03421
B	244	GLU	-	expression tag	UNP P03421
B	245	ASN	-	expression tag	UNP P03421
B	246	LEU	-	expression tag	UNP P03421
B	247	TYR	-	expression tag	UNP P03421
B	248	PHE	-	expression tag	UNP P03421
B	249	GLN	-	expression tag	UNP P03421
B	250	GLY	-	expression tag	UNP P03421
B	251	HIS	-	expression tag	UNP P03421
B	252	HIS	-	expression tag	UNP P03421
B	253	HIS	-	expression tag	UNP P03421
B	254	HIS	-	expression tag	UNP P03421
B	255	HIS	-	expression tag	UNP P03421
B	256	HIS	-	expression tag	UNP P03421
C	171	VAL	ILE	variant	UNP P03421
C	242	LYS	-	expression tag	UNP P03421

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Chain	Residue	Modelled	Actual	Comment	Reference
C	243	GLY	-	expression tag	UNP P03421
C	244	GLU	-	expression tag	UNP P03421
C	245	ASN	-	expression tag	UNP P03421
C	246	LEU	-	expression tag	UNP P03421
C	247	TYR	-	expression tag	UNP P03421
C	248	PHE	-	expression tag	UNP P03421
C	249	GLN	-	expression tag	UNP P03421
C	250	GLY	-	expression tag	UNP P03421
C	251	HIS	-	expression tag	UNP P03421
C	252	HIS	-	expression tag	UNP P03421
C	253	HIS	-	expression tag	UNP P03421
C	254	HIS	-	expression tag	UNP P03421
C	255	HIS	-	expression tag	UNP P03421
C	256	HIS	-	expression tag	UNP P03421
D	171	VAL	ILE	variant	UNP P03421
D	242	LYS	-	expression tag	UNP P03421
D	243	GLY	-	expression tag	UNP P03421
D	244	GLU	-	expression tag	UNP P03421
D	245	ASN	-	expression tag	UNP P03421
D	246	LEU	-	expression tag	UNP P03421
D	247	TYR	-	expression tag	UNP P03421
D	248	PHE	-	expression tag	UNP P03421
D	249	GLN	-	expression tag	UNP P03421
D	250	GLY	-	expression tag	UNP P03421
D	251	HIS	-	expression tag	UNP P03421
D	252	HIS	-	expression tag	UNP P03421
D	253	HIS	-	expression tag	UNP P03421
D	254	HIS	-	expression tag	UNP P03421
D	255	HIS	-	expression tag	UNP P03421
D	256	HIS	-	expression tag	UNP P03421
E	171	VAL	ILE	variant	UNP P03421
E	242	LYS	-	expression tag	UNP P03421
E	243	GLY	-	expression tag	UNP P03421
E	244	GLU	-	expression tag	UNP P03421
E	245	ASN	-	expression tag	UNP P03421
E	246	LEU	-	expression tag	UNP P03421
E	247	TYR	-	expression tag	UNP P03421
E	248	PHE	-	expression tag	UNP P03421
E	249	GLN	-	expression tag	UNP P03421
E	250	GLY	-	expression tag	UNP P03421
E	251	HIS	-	expression tag	UNP P03421
E	252	HIS	-	expression tag	UNP P03421

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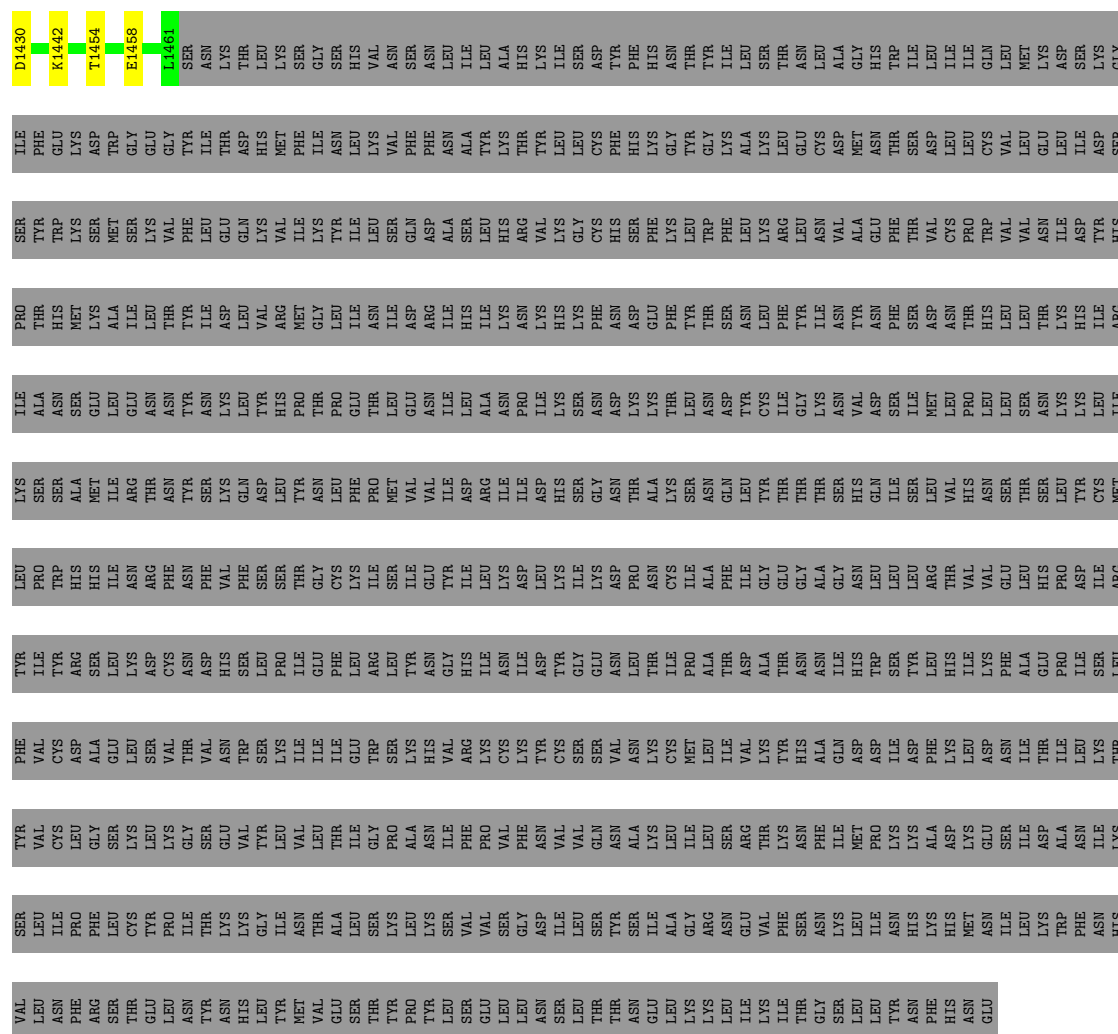
Chain	Residue	Modelled	Actual	Comment	Reference
E	253	HIS	-	expression tag	UNP P03421
E	254	HIS	-	expression tag	UNP P03421
E	255	HIS	-	expression tag	UNP P03421
E	256	HIS	-	expression tag	UNP P03421

- Molecule 3 is a protein called Compound 16.

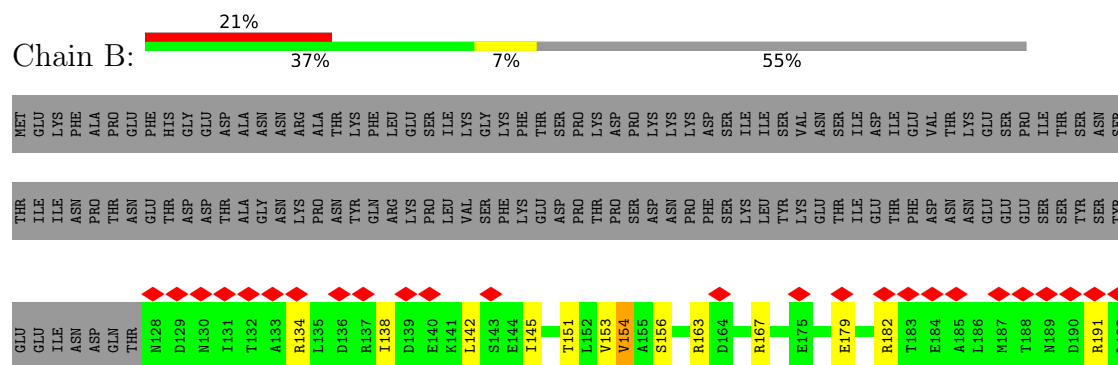
Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	3	Total	C	Cl	N	O	0	0
			36	28	1	4	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		AltConf
4	A	2	Total	O	0
			2	2	
4	G	2	Total	O	0
			2	2	

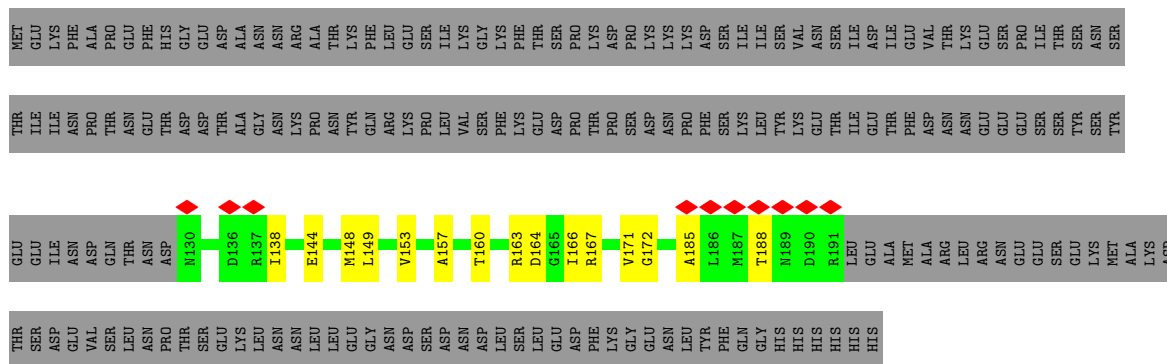


- Molecule 2: Phosphoprotein

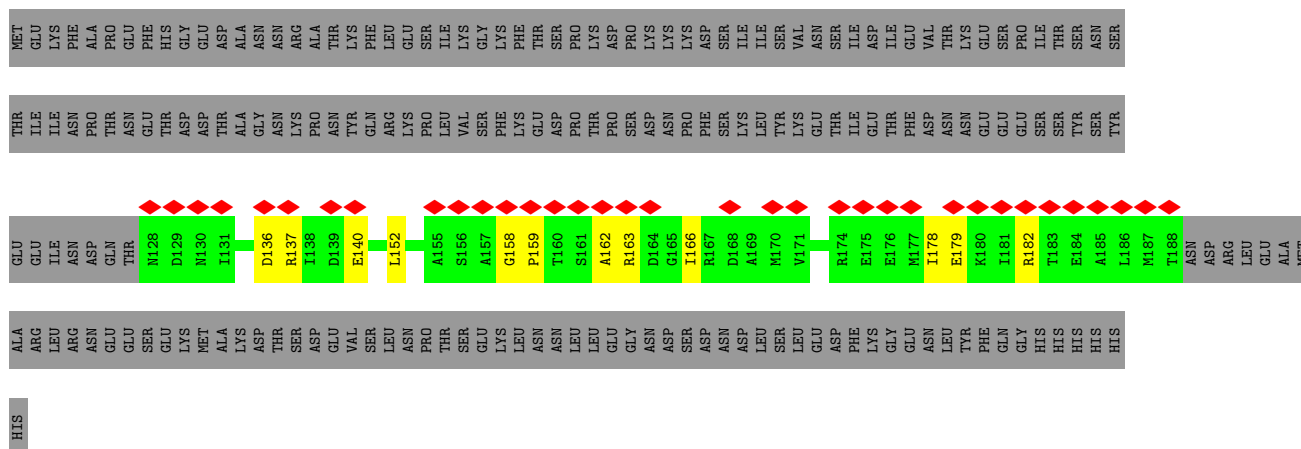


- Molecule 2: Phosphoprotein

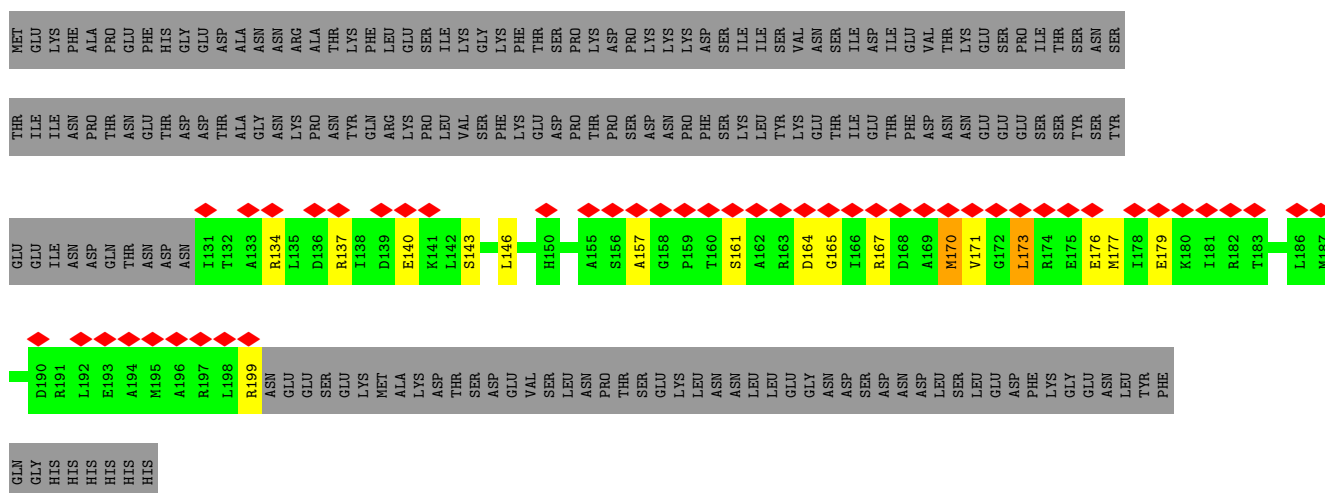




- Molecule 2: Phosphoprotein



- Molecule 2: Phosphoprotein



- Molecule 3: Compound 16





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	1127112	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS GLACIOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	2.771	Depositor
Minimum map value	-2.064	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.072	Depositor
Recommended contour level	0.378	Depositor
Map size (Å)	273.02402, 273.02402, 273.02402	wwPDB
Map dimensions	288, 288, 288	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.9480001, 0.9480001, 0.9480001	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: A1BIT, A1BIY, HPE

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	0.32	0/11387	0.43	0/15414
2	B	0.86	2/888 (0.2%)	1.43	6/1193 (0.5%)
2	C	0.34	0/479	0.55	1/642 (0.2%)
2	D	0.31	0/468	0.54	0/628
2	E	0.32	0/536	0.63	1/717 (0.1%)
All	All	0.38	2/13758 (0.0%)	0.57	8/18594 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	154	VAL	C-N	-17.87	1.07	1.33
2	B	153	VAL	C-N	-16.67	1.10	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	154	VAL	O-C-N	-29.62	84.08	122.88
2	B	154	VAL	CA-C-N	17.12	147.43	122.09
2	B	154	VAL	C-N-CA	17.12	147.43	122.09
2	B	153	VAL	CA-C-N	15.77	144.55	123.11
2	B	153	VAL	C-N-CA	15.77	144.55	123.11
2	B	153	VAL	O-C-N	-14.97	107.87	123.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	172	GLY	CA-C-O	-7.13	117.58	122.22
2	E	171	VAL	N-CA-C	-5.31	108.11	113.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	154	VAL	Mainchain

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	11153	0	11354	110	0
2	B	885	0	870	13	0
2	C	478	0	496	14	0
2	D	467	0	483	13	0
2	E	535	0	563	12	0
3	G	36	0	10	1	0
4	A	2	0	0	1	0
4	G	2	0	0	0	0
All	All	13558	0	13776	144	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 5.

All (144) 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)
4:A:2201:HOH:O	3:G:3:A1BIY:CL34	2.35	0.81
1:A:489:LEU:O	2:D:137:ARG:NH2	2.17	0.75
1:A:403:LYS:NZ	2:B:236:LEU:O	2.23	0.72
1:A:189:LEU:HD11	1:A:299:LEU:HD11	1.72	0.71
1:A:869:MET:HG3	1:A:872:THR:HB	1.71	0.71
1:A:1037:ASN:ND2	1:A:1192:LEU:O	2.24	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1357:ARG:NH2	1:A:1400:GLU:OE1	2.27	0.68
1:A:62:LYS:O	1:A:68:GLN:NE2	2.28	0.66
1:A:555:ARG:O	1:A:563:GLN:NE2	2.28	0.66
1:A:1303:ASP:OD1	1:A:1303:ASP:N	2.28	0.64
1:A:468:ARG:NH2	1:A:657:GLU:OE2	2.24	0.64
1:A:595:CYS:O	1:A:599:ASN:ND2	2.31	0.64
2:E:157:ALA:O	2:E:161:SER:OG	2.16	0.63
1:A:995:ARG:NH1	1:A:999:ARG:O	2.31	0.63
1:A:1335:ASN:ND2	1:A:1380:ASP:OD2	2.32	0.62
1:A:468:ARG:NH1	1:A:794:ASP:OD2	2.33	0.62
1:A:901:PHE:HZ	1:A:937:ILE:HG12	1.65	0.61
2:B:145:ILE:HG12	2:E:146:LEU:HD13	1.82	0.61
1:A:1224:ILE:HG13	1:A:1387:ASN:HB3	1.83	0.61
1:A:514:ILE:HD13	2:C:148:MET:HE2	1.83	0.61
1:A:1204:LYS:NZ	1:A:1208:GLU:OE2	2.33	0.60
1:A:101:GLU:OE2	1:A:213:LYS:NZ	2.29	0.59
2:C:160:THR:HG23	2:C:163:ARG:HH21	1.66	0.59
1:A:1171:LYS:O	1:A:1178:ARG:NH2	2.32	0.59
1:A:765:GLU:HG2	1:A:770:TYR:HA	1.83	0.58
1:A:874:GLN:HE21	1:A:877:GLY:HA2	1.69	0.58
2:D:158:GLY:H	2:D:159:PRO:HD2	1.69	0.57
1:A:1046:PHE:O	1:A:1209:ARG:NH2	2.27	0.57
1:A:115:ARG:O	1:A:119:ILE:HG12	2.03	0.57
1:A:1003:ASP:OD1	1:A:1207:ARG:NH2	2.38	0.57
2:E:176:GLU:O	2:E:179:GLU:HG3	2.04	0.57
1:A:1189:ARG:NH1	1:A:1190:GLU:O	2.38	0.57
1:A:437:ARG:NH2	1:A:750:PRO:O	2.39	0.56
1:A:1312:LEU:O	1:A:1316:THR:OG1	2.20	0.56
1:A:509:THR:HG22	1:A:510:GLU:H	1.71	0.56
2:B:191:ARG:HG2	2:B:195:MET:HE2	1.88	0.56
2:C:153:VAL:HG22	2:D:152:LEU:HG	1.88	0.56
1:A:47:ASN:OD1	1:A:48:LEU:N	2.39	0.55
1:A:1059:ARG:NH2	1:A:1104:GLU:OE2	2.21	0.55
2:C:188:THR:HG22	2:D:182:ARG:HE	1.71	0.55
1:A:796:ILE:HG12	1:A:847:GLU:HG3	1.89	0.55
1:A:1064:LEU:O	1:A:1210:SER:OG	2.09	0.55
2:B:142:LEU:HD11	2:C:138:ILE:HG23	1.89	0.54
1:A:1430:ASP:OD1	1:A:1430:ASP:N	2.40	0.54
1:A:544:PRO:HG2	1:A:549:ILE:HD12	1.88	0.54
1:A:1351:ALA:HB3	1:A:1353:ILE:HG22	1.90	0.53
1:A:1296:ASP:OD1	1:A:1305:LYS:NZ	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1110:ILE:HG23	1:A:1111:MET:HG2	1.89	0.53
1:A:1240:GLY:HA3	1:A:1412:LEU:O	2.08	0.53
2:D:162:ALA:O	2:D:166:ILE:HG13	2.09	0.53
1:A:648:ALA:HB2	1:A:783:LYS:HE2	1.91	0.53
2:B:156:SER:O	2:B:167:ARG:NH1	2.39	0.52
1:A:851:ILE:HG13	1:A:853:HIS:HD2	1.74	0.52
1:A:186:LYS:NZ	1:A:1018:TYR:O	2.43	0.52
1:A:816:ASP:OD1	1:A:867:GLN:NE2	2.43	0.52
2:C:164:ASP:OD2	2:D:163:ARG:NH1	2.43	0.52
1:A:461:LEU:HD23	1:A:715:ILE:HD11	1.93	0.51
1:A:1259:THR:HG22	1:A:1260:LYS:H	1.75	0.51
1:A:1007:GLU:OE1	1:A:1223:SER:OG	2.28	0.49
1:A:897:ILE:HD12	1:A:1145:LYS:HG3	1.94	0.49
2:E:140:GLU:O	2:E:143:SER:HB3	2.13	0.49
2:C:167:ARG:O	2:C:171:VAL:HG22	2.13	0.48
1:A:734:HIS:CE1	1:A:775:GLY:HA3	2.48	0.48
1:A:554:PRO:HD3	1:A:747:ARG:HG3	1.95	0.48
1:A:952:TYR:O	1:A:955:ILE:HG22	2.12	0.48
1:A:1370:ARG:O	1:A:1374:GLU:HG2	2.13	0.48
1:A:319:CYS:SG	2:B:238:LEU:HD22	2.54	0.48
2:B:134:ARG:O	2:B:138:ILE:HG12	2.13	0.48
1:A:526:ARG:NH2	1:A:1458:GLU:OE1	2.47	0.48
1:A:1353:ILE:HD11	1:A:1357:ARG:HG2	1.96	0.48
1:A:372:GLN:HG3	1:A:392:ILE:HG23	1.96	0.47
1:A:448:ASN:HD21	2:B:163:ARG:HH22	1.62	0.47
1:A:36:PHE:CE2	1:A:81:LEU:HD21	2.49	0.47
2:B:179:GLU:OE1	2:B:182:ARG:NH2	2.31	0.47
1:A:565:TYR:CZ	1:A:569:GLU:HG3	2.50	0.47
2:C:149:LEU:O	2:C:153:VAL:HG23	2.14	0.46
1:A:348:GLU:OE2	1:A:863:SER:OG	2.34	0.46
2:D:158:GLY:H	2:D:159:PRO:CD	2.29	0.46
2:E:137:ARG:HA	2:E:140:GLU:HG2	1.96	0.46
1:A:576:SER:OG	1:A:585:TYR:O	2.33	0.46
2:E:170:MET:HB3	2:E:170:MET:HE3	1.56	0.46
1:A:804:ILE:HG23	1:A:815:ILE:HG23	1.97	0.46
1:A:983:MET:HG3	1:A:989:ASP:O	2.16	0.45
1:A:403:LYS:HE3	2:B:234:ASN:HB3	1.98	0.45
1:A:193:LEU:HD11	1:A:197:MET:HE3	1.98	0.45
1:A:423:TYR:CG	1:A:871:LYS:HE2	2.51	0.45
1:A:462:ARG:NH1	1:A:517:SER:O	2.48	0.45
1:A:39:PRO:HA	1:A:47:ASN:HD22	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:601:VAL:HG21	1:A:642:LEU:HD23	1.98	0.45
1:A:1454:THR:HG23	2:E:199:ARG:HD2	1.98	0.44
1:A:535:MET:HE3	1:A:641:ILE:HG22	1.99	0.44
1:A:912:THR:HG21	1:A:929:ARG:HB2	1.98	0.44
1:A:1266:SER:HA	1:A:1386:GLN:NE2	2.32	0.44
2:B:216:LEU:HD22	2:B:220:SER:OG	2.16	0.44
2:D:136:ASP:O	2:D:140:GLU:HG3	2.17	0.44
1:A:872:THR:HG23	1:A:879:TYR:HE1	1.82	0.44
1:A:1240:GLY:O	1:A:1367:PRO:HG2	2.18	0.44
1:A:866:MET:HG2	1:A:867:GLN:N	2.33	0.44
1:A:1031:LEU:O	1:A:1033:ASP:N	2.50	0.44
1:A:1106:ASP:OD2	1:A:1442:LYS:NZ	2.45	0.44
1:A:1200:THR:OG1	1:A:1418:GLU:OE1	2.35	0.44
1:A:494:LEU:O	1:A:498:LYS:HG2	2.17	0.43
1:A:1042:CYS:SG	1:A:1184:CYS:HA	2.57	0.43
1:A:440:MET:HE2	1:A:440:MET:HB3	1.71	0.43
1:A:1077:ASN:O	1:A:1081:VAL:HG23	2.19	0.43
2:C:166:ILE:HD13	2:D:163:ARG:HD2	2.00	0.43
1:A:1127:GLU:OE2	1:A:1343:SER:OG	2.35	0.43
1:A:242:PHE:CE2	1:A:255:HIS:HD2	2.37	0.42
1:A:99:SER:O	1:A:103:ILE:HG12	2.19	0.42
1:A:73:LYS:HE3	1:A:73:LYS:HB2	1.71	0.42
2:C:160:THR:HG23	2:C:163:ARG:NH2	2.33	0.42
2:D:179:GLU:OE1	2:D:182:ARG:NH1	2.53	0.42
1:A:755:ASP:O	1:A:756:HIS:HB2	2.19	0.42
2:D:158:GLY:N	2:D:159:PRO:HD2	2.34	0.42
1:A:550:TRP:HB2	1:A:566:ILE:HD11	2.02	0.42
1:A:950:LYS:HD3	1:A:950:LYS:HA	1.73	0.42
2:E:134:ARG:HG3	2:E:137:ARG:HH21	1.85	0.41
1:A:209:ASN:O	1:A:213:LYS:HG2	2.20	0.41
1:A:473:PHE:CD1	1:A:481:PRO:HG3	2.55	0.41
1:A:698:ILE:O	1:A:860:THR:HA	2.19	0.41
1:A:1133:LYS:HE2	1:A:1133:LYS:HB3	1.77	0.41
2:E:164:ASP:OD1	2:E:165:GLY:N	2.53	0.41
1:A:1230:ILE:HG12	1:A:1395:LEU:HD21	2.02	0.41
1:A:1102:THR:HG22	1:A:1442:LYS:NZ	2.36	0.41
2:E:177:MET:HE3	2:E:177:MET:HB2	1.99	0.41
1:A:456:SER:OG	2:C:148:MET:HE1	2.21	0.41
1:A:710:TYR:OH	2:C:157:ALA:O	2.37	0.41
1:A:62:LYS:HA	1:A:67:THR:HG21	2.03	0.41
1:A:901:PHE:CZ	1:A:937:ILE:HG12	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:THR:HG23	1:A:246:LEU:O	2.21	0.41
1:A:911:LEU:O	1:A:914:GLU:HG2	2.21	0.41
1:A:970:ASN:OD1	1:A:970:ASN:N	2.51	0.41
1:A:1116:PRO:HG2	1:A:1427:PHE:CD1	2.55	0.41
2:E:173:LEU:HD13	2:E:173:LEU:HA	1.84	0.41
1:A:355:ARG:NH1	2:B:211:SER:O	2.53	0.40
2:B:151:THR:O	2:B:151:THR:HG22	2.22	0.40
1:A:64:LEU:O	1:A:68:GLN:HG2	2.22	0.40
1:A:435:ASP:HB3	1:A:438:GLN:HG2	2.03	0.40
1:A:902:LYS:HE2	1:A:902:LYS:HB3	1.91	0.40
2:C:185:ALA:HB2	2:D:178:ILE:HG23	2.03	0.40
1:A:511:ARG:HH11	2:C:144:GLU:CD	2.27	0.40
1:A:29:ASN:O	1:A:33:SER:OG	2.24	0.40
1:A:482:THR:HG22	1:A:522:TYR:HD2	1.86	0.40
2:D:159:PRO:HG3	2:E:167:ARG:HB2	2.03	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	1359/2201 (62%)	1288 (95%)	66 (5%)	5 (0%)	30	60
2	B	112/256 (44%)	108 (96%)	4 (4%)	0	100	100
2	C	60/256 (23%)	56 (93%)	4 (7%)	0	100	100
2	D	59/256 (23%)	56 (95%)	3 (5%)	0	100	100
2	E	67/256 (26%)	58 (87%)	9 (13%)	0	100	100
All	All	1657/3225 (51%)	1566 (94%)	86 (5%)	5 (0%)	37	66

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	490	PRO
1	A	325	HIS
1	A	326	ASN
1	A	892	PRO
1	A	949	ASN

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	1270/2041 (62%)	1268 (100%)	2 (0%)	87	96
2	B	100/233 (43%)	100 (100%)	0	100	100
2	C	52/233 (22%)	52 (100%)	0	100	100
2	D	51/233 (22%)	51 (100%)	0	100	100
2	E	57/233 (24%)	55 (96%)	2 (4%)	32	67
All	All	1530/2973 (52%)	1526 (100%)	4 (0%)	84	95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1055	VAL
1	A	1200	THR
2	E	170	MET
2	E	173	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	HIS
1	A	255	HIS
1	A	312	GLN
1	A	417	ASN
1	A	448	ASN
1	A	478	ASN

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Mol	Chain	Res	Type
1	A	603	ASN
1	A	705	ASN
1	A	748	HIS
1	A	827	GLN
1	A	874	GLN
1	A	930	ASN
1	A	1022	HIS
1	A	1185	ASN
1	A	1249	ASN
1	A	1362	HIS
1	A	1369	ASN
1	A	1438	GLN
1	A	1441	GLN
2	C	130	ASN
2	C	189	ASN
2	E	189	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	A1BIY	G	3	3	11,13,13	2.65	2 (18%)	13,17,17	1.44	3 (23%)
3	A1BIT	G	1	3	9,11,12	1.13	1 (11%)	8,14,16	1.50	2 (25%)
3	HPE	G	2	3	11,12,13	0.73	0	9,14,16	0.78	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	A1BIY	G	3	3	-	0/10/10/10	0/1/1/1
3	A1BIT	G	1	3	-	1/5/17/19	0/1/1/1
3	HPE	G	2	3	-	0/6/7/9	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	3	A1BIY	C03-N02	7.98	1.44	1.33
3	G	3	A1BIY	O04-C03	-2.77	1.18	1.23
3	G	1	A1BIT	C5-C4	-2.30	1.49	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	3	A1BIY	C05-C03-N02	3.26	118.74	114.62
3	G	1	A1BIT	C5-C4-C2	-3.15	108.25	113.60
3	G	1	A1BIT	O1-C9-C8	-2.42	118.45	124.86
3	G	3	A1BIY	O04-C03-N02	-2.01	119.77	123.13
3	G	3	A1BIY	C01-N02-C03	-2.01	118.71	122.21

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	G	1	A1BIT	C4-C8-C9-O1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	3	A1BIY	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	B	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	153:VAL	C	154:VAL	N	1.10
1	B	154:VAL	C	155:ALA	N	1.07

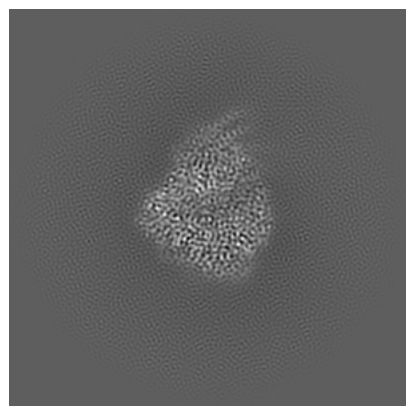
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-47927. These allow visual inspection of the internal detail of the map and identification of artifacts.

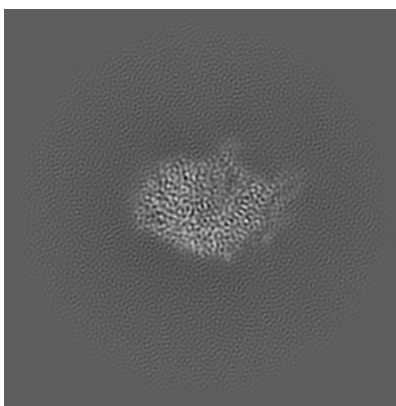
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

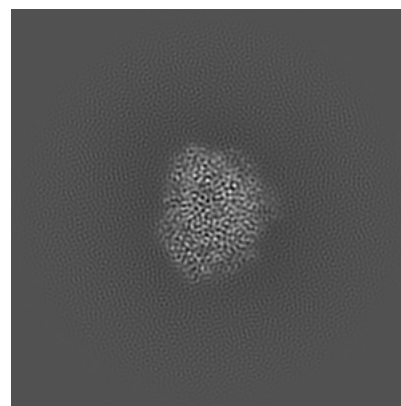
6.1.1 Primary map



X

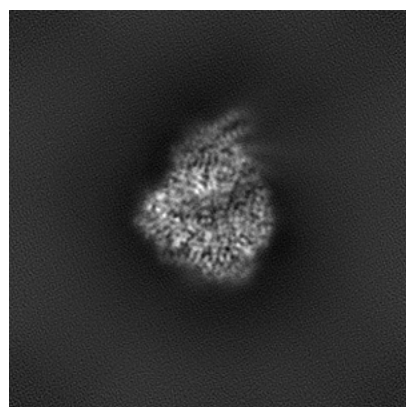


Y

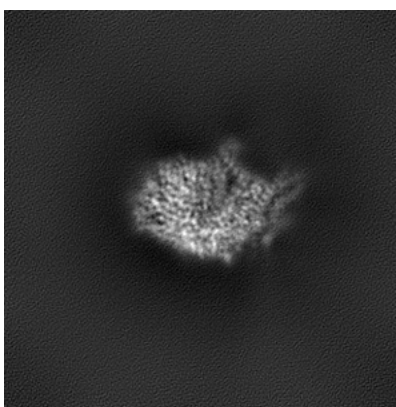


Z

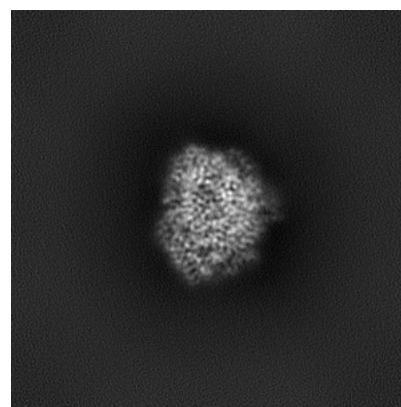
6.1.2 Raw map



X



Y

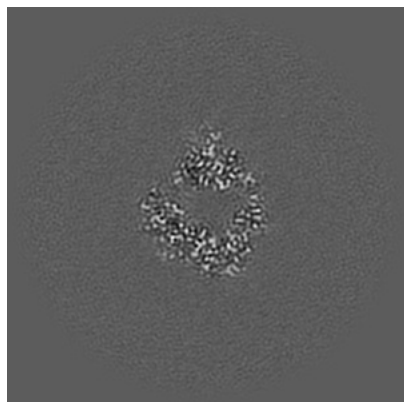


Z

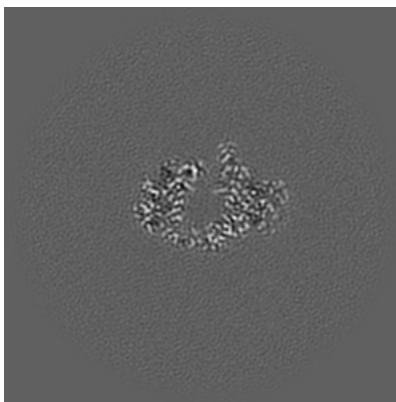
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

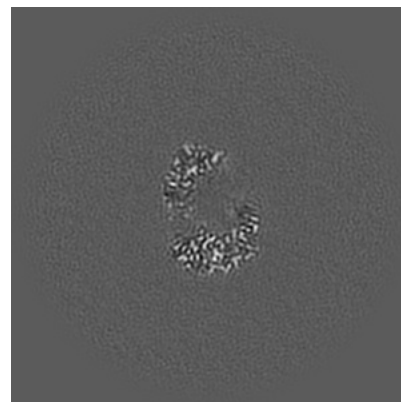
6.2.1 Primary map



X Index: 144

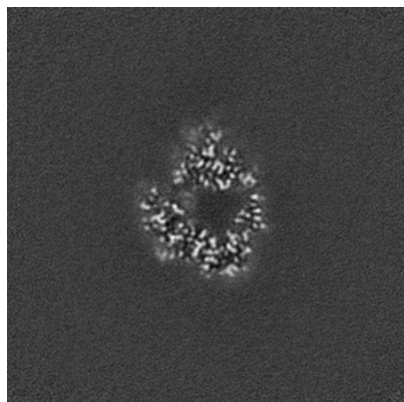


Y Index: 144

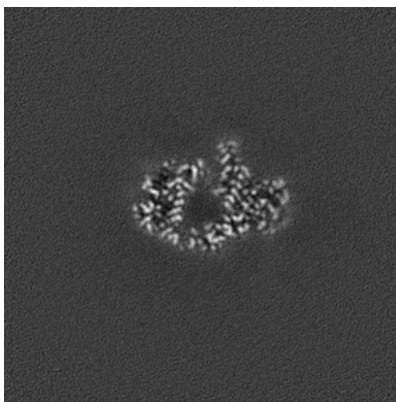


Z Index: 144

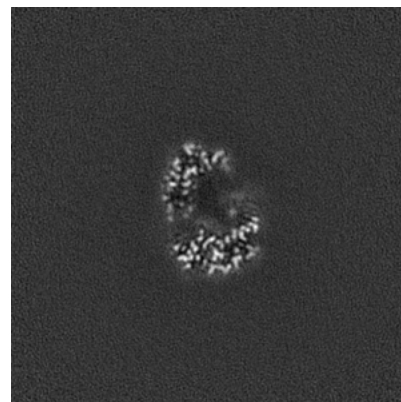
6.2.2 Raw map



X Index: 144



Y Index: 144

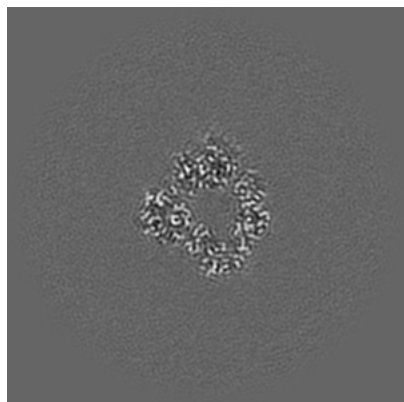


Z Index: 144

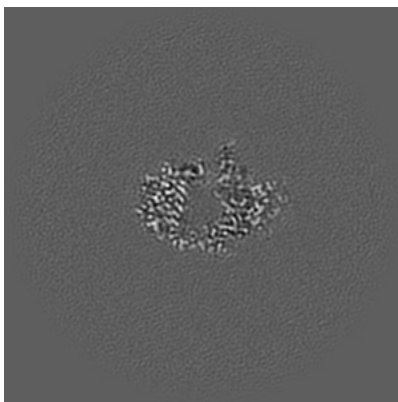
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

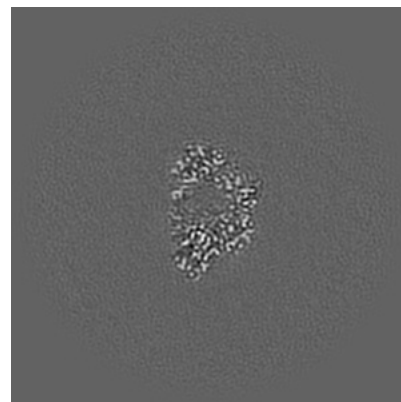
6.3.1 Primary map



X Index: 137

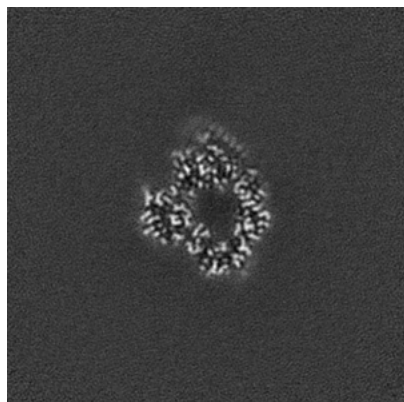


Y Index: 142

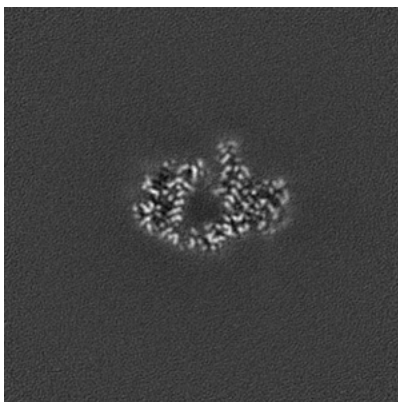


Z Index: 132

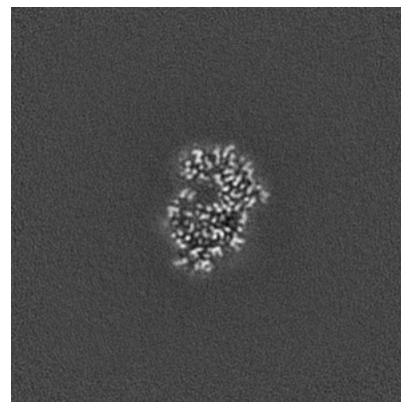
6.3.2 Raw map



X Index: 137



Y Index: 144

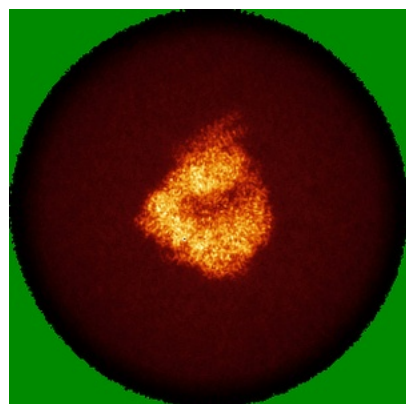


Z Index: 127

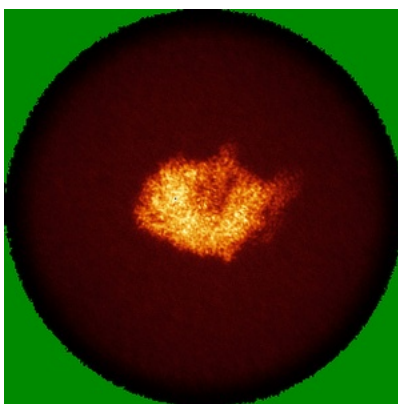
The images above show the largest variance slices of the map in three orthogonal directions.

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

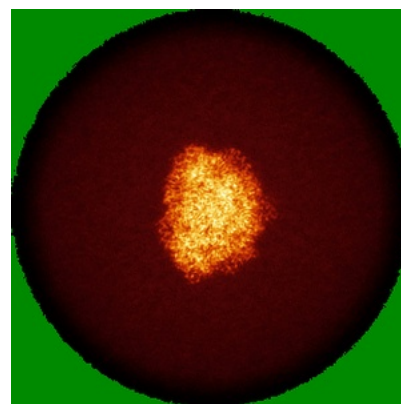
6.4.1 Primary map



X

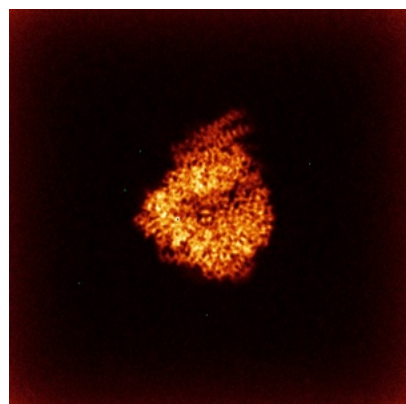


Y

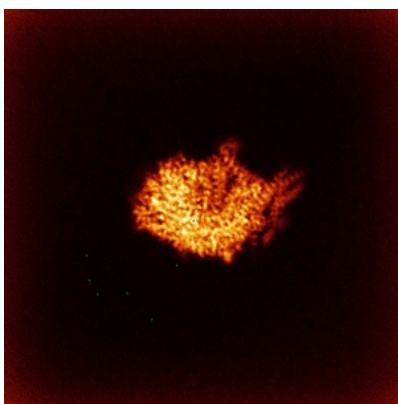


Z

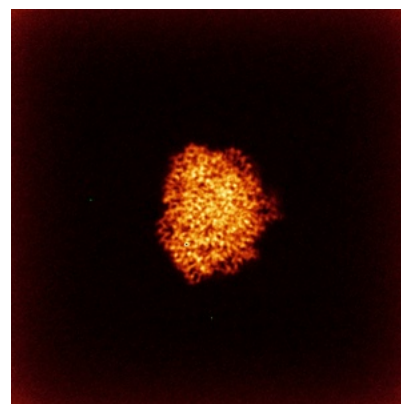
6.4.2 Raw map



X



Y

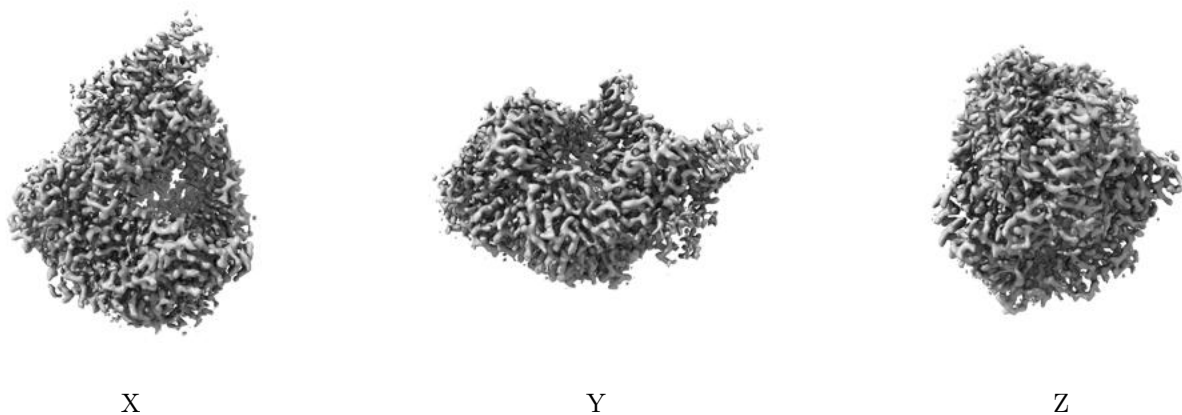


Z

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.

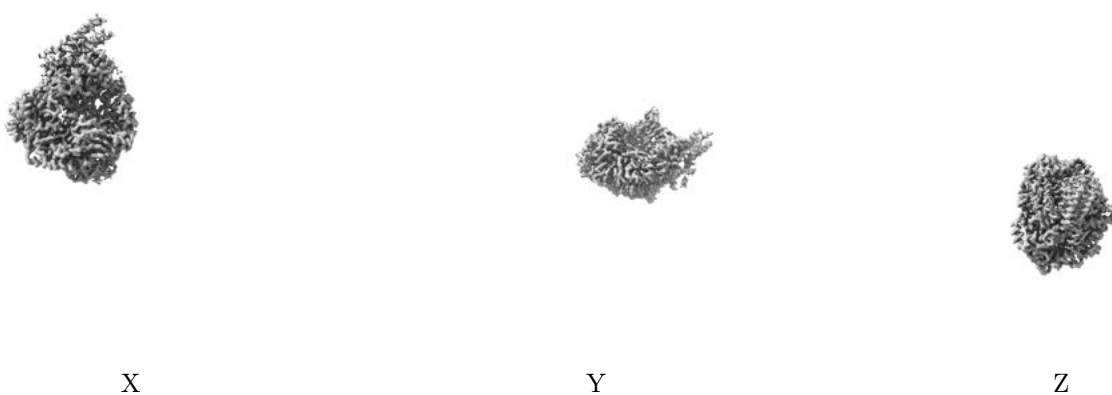
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.378. 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.

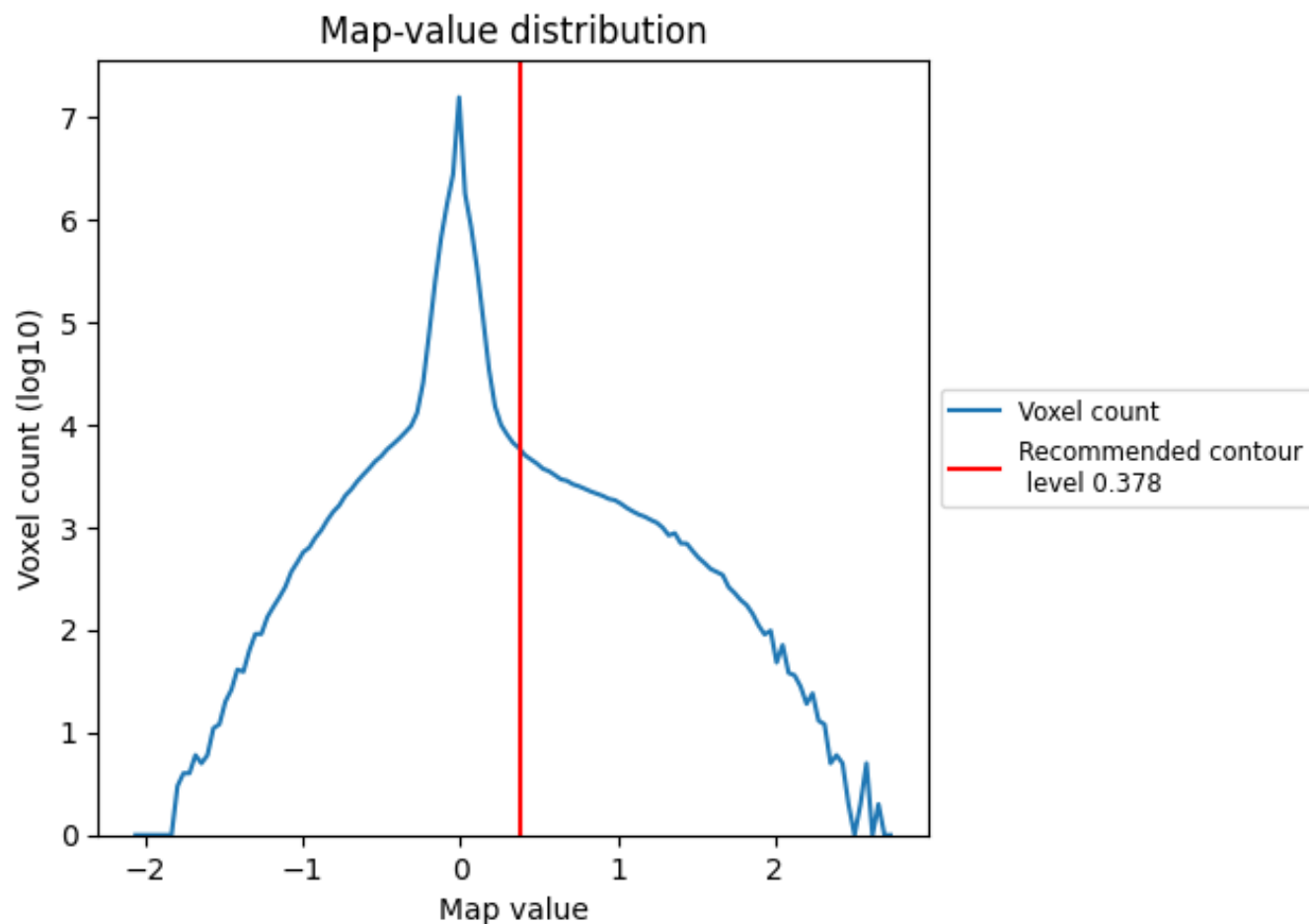
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

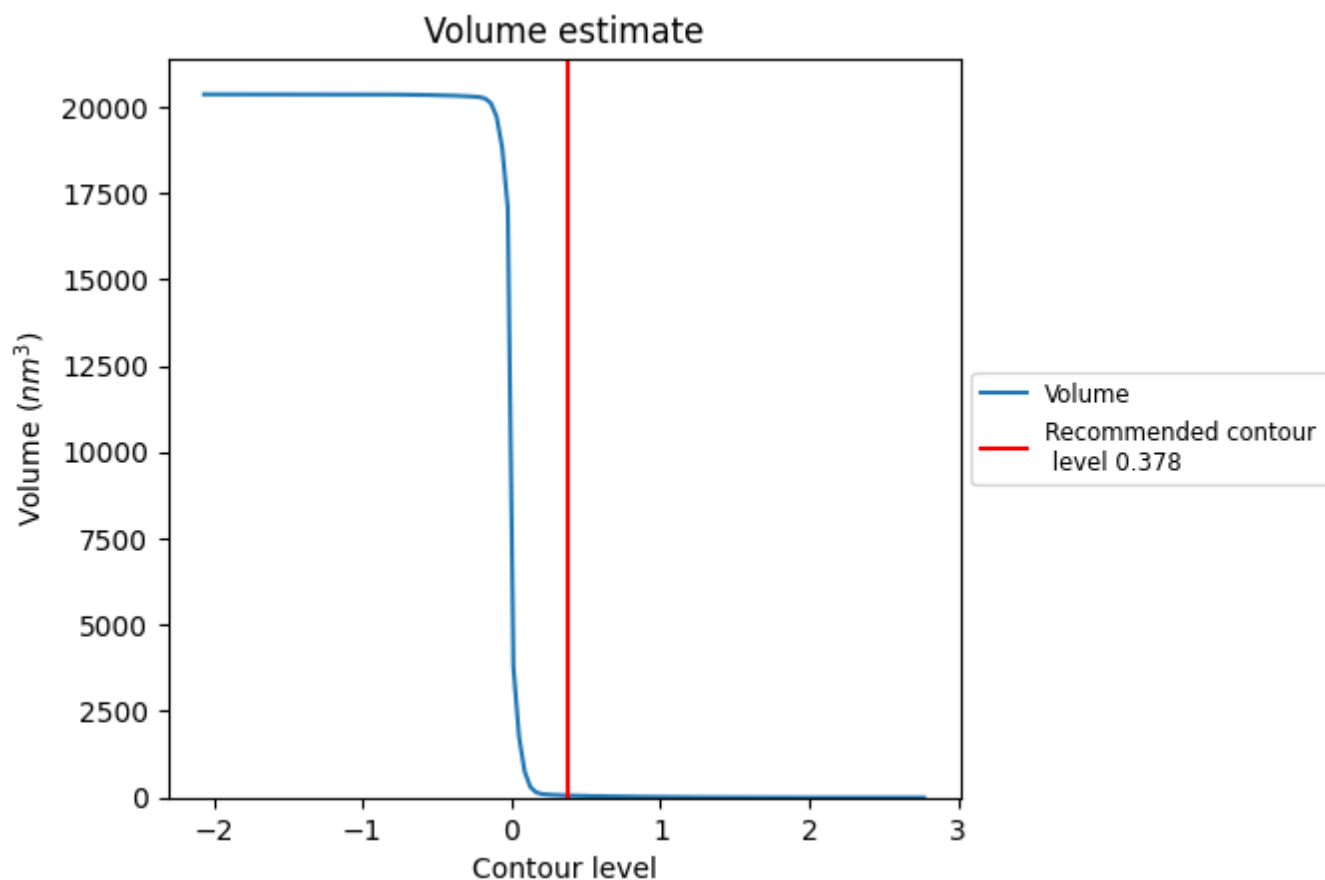
This section contains the results of statistical analysis of the map.

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.

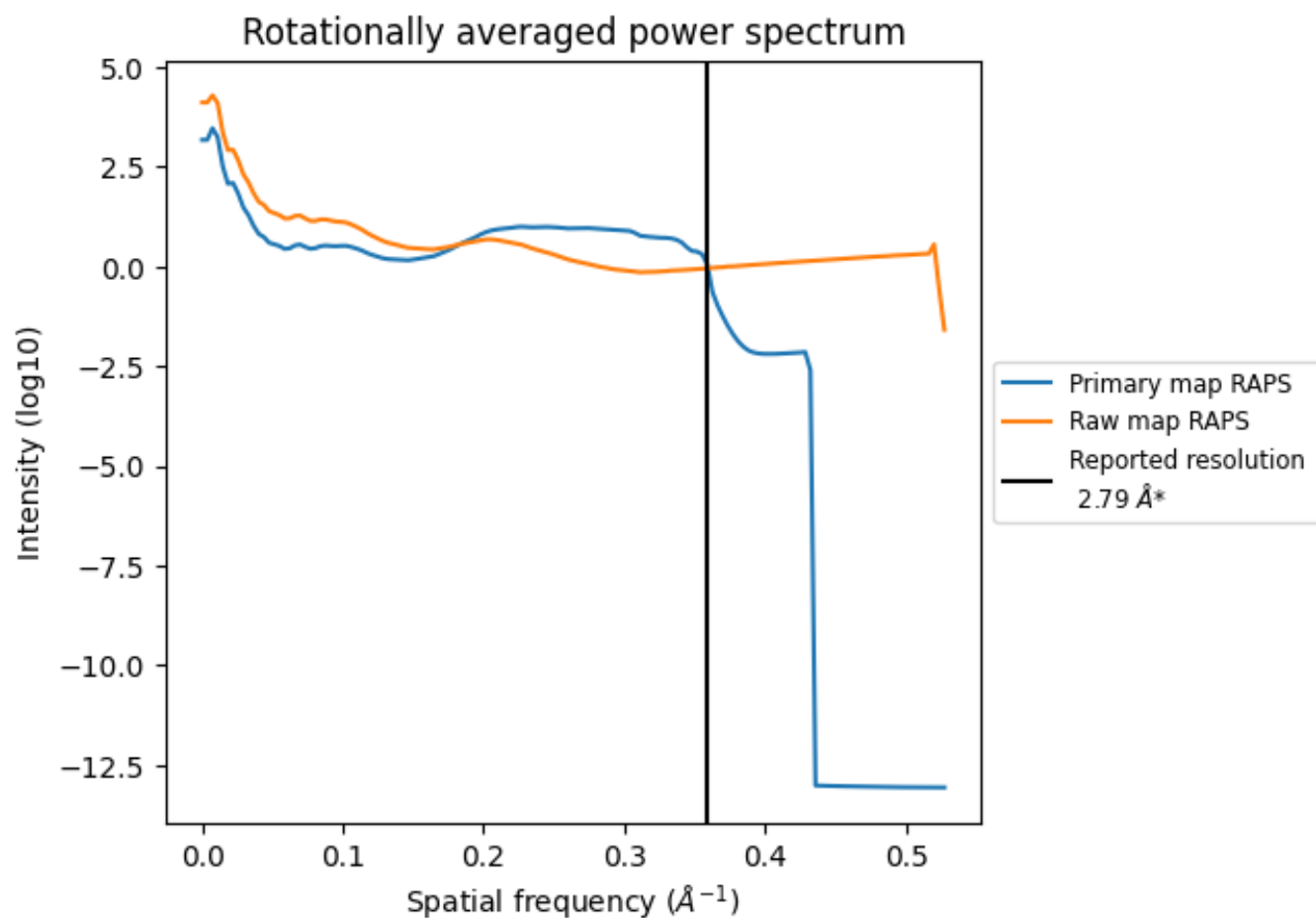
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 60 nm³; this corresponds to an approximate mass of 54 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 ⓘ

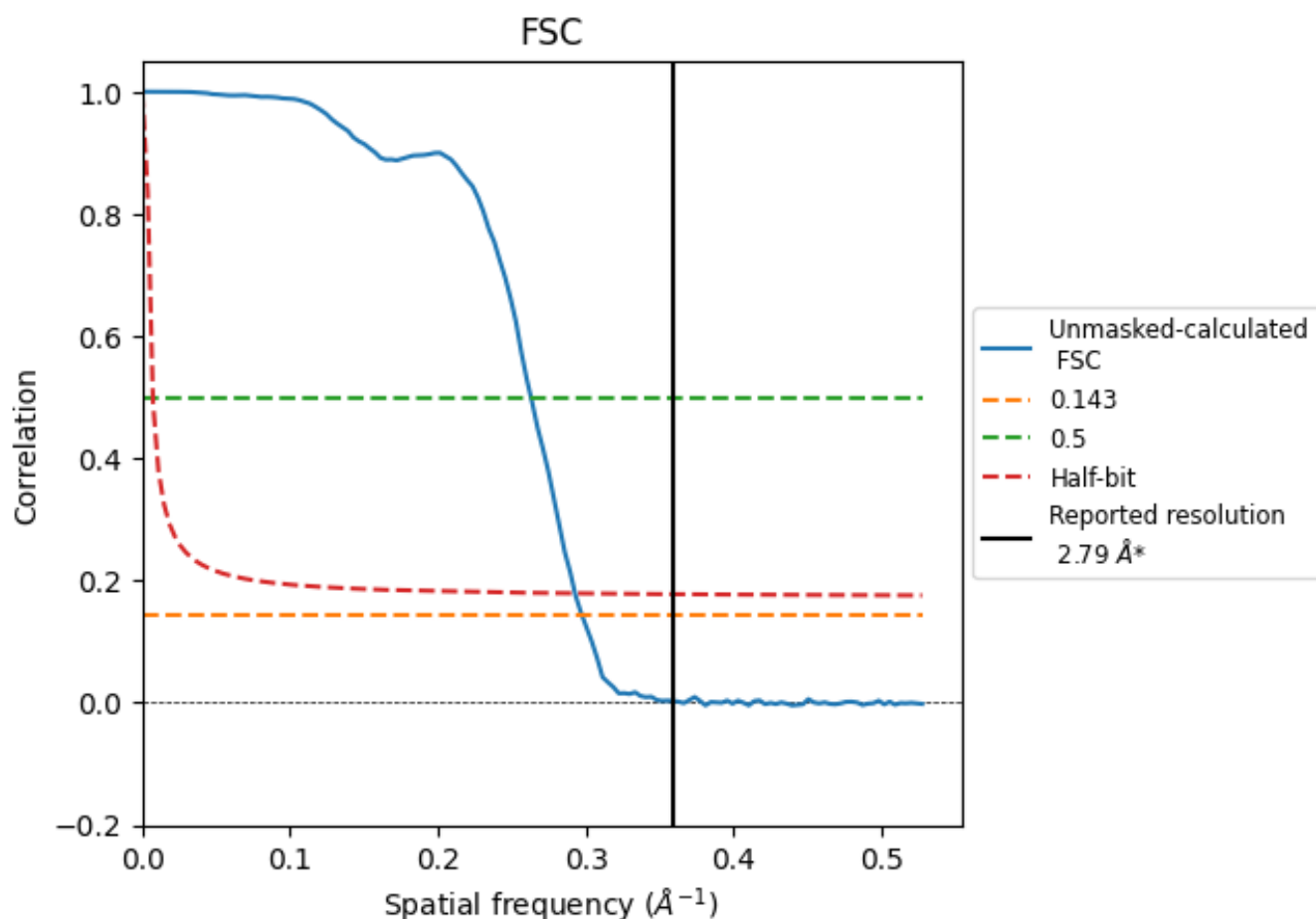


*Reported resolution corresponds to spatial frequency of 0.358 \AA^{-1}

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 0.358 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.79	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.36	3.80	3.42

*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 3.36 differs from the reported value 2.79 by more than 10 %

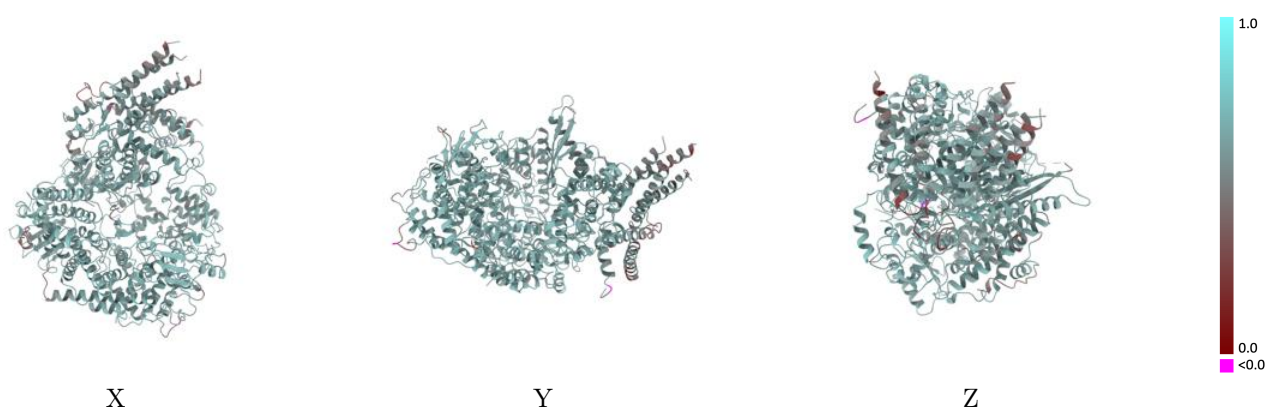
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-47927 and PDB model 9ECV. Per-residue inclusion information can be found in section 3 on page 8.

9.1 Map-model overlay [i](#)

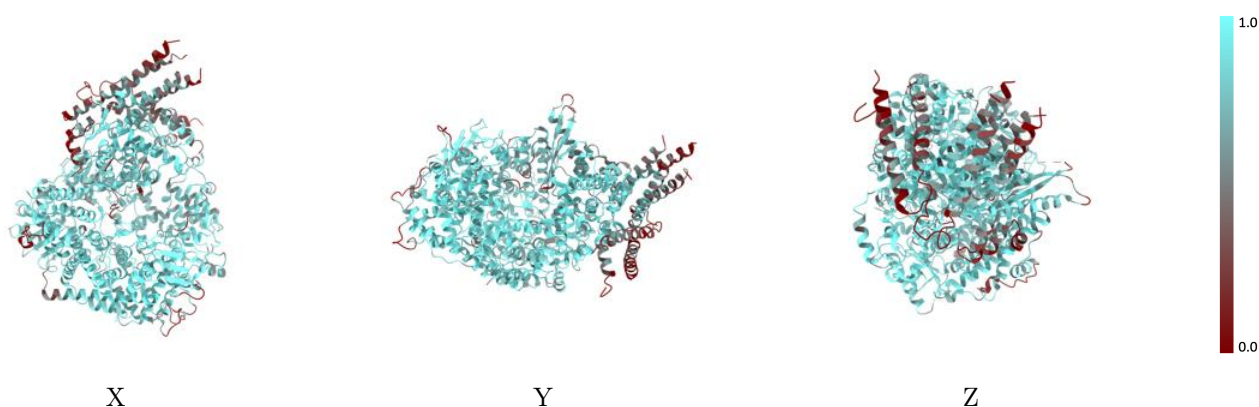
This section was not generated.

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



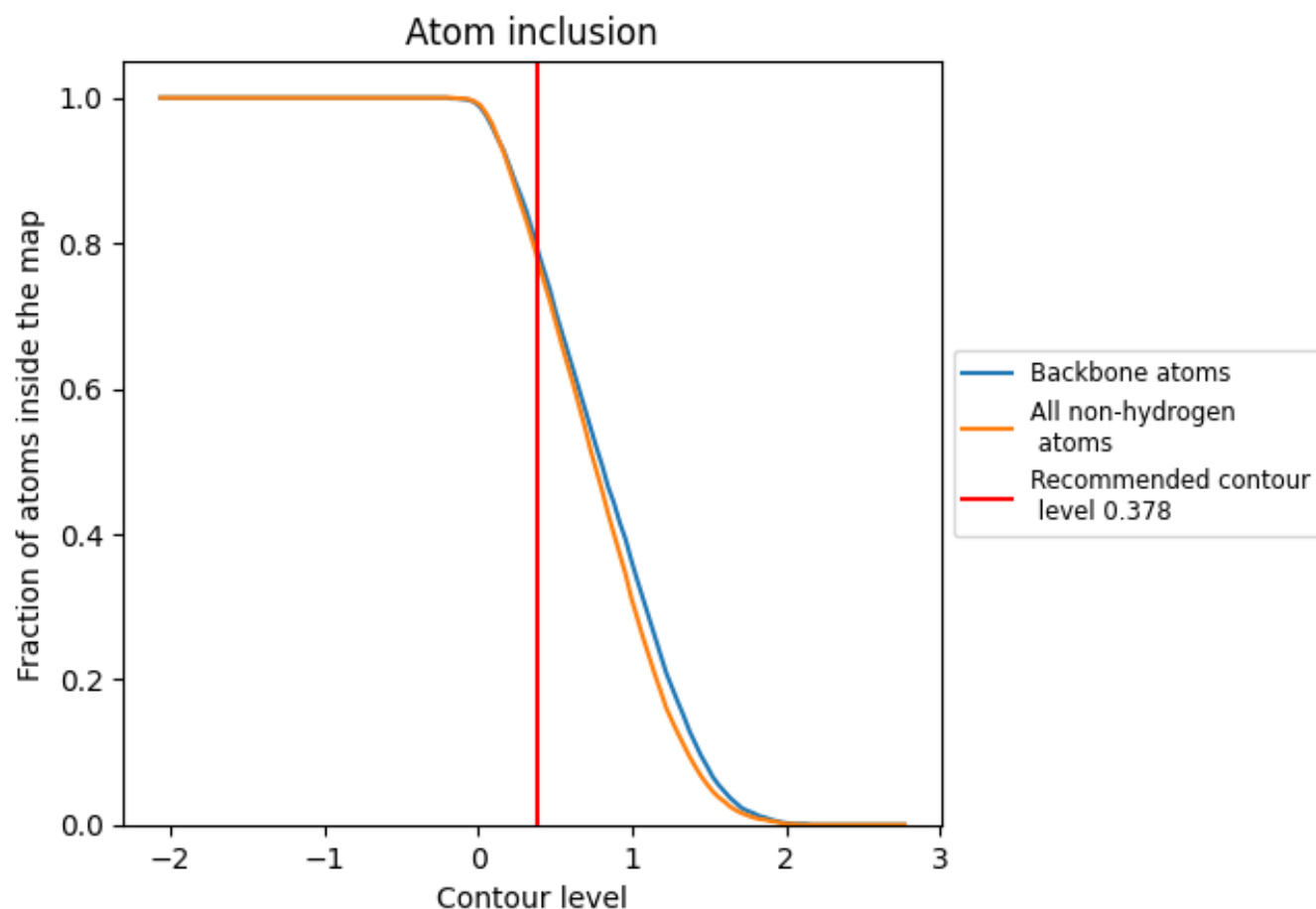
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.378).

9.4 Atom inclusion [i](#)



At the recommended contour level, 80% of all backbone atoms, 78% 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 (0.378) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div>0.7850</div>	<div><div></div>0.6080</div>
A	<div><div></div>0.8590</div>	<div><div></div>0.6250</div>
B	<div><div></div>0.4350</div>	<div><div></div>0.5320</div>
C	<div><div></div>0.6770</div>	<div><div></div>0.5820</div>
D	<div><div></div>0.3800</div>	<div><div></div>0.5080</div>
E	<div><div></div>0.2940</div>	<div><div></div>0.4850</div>
G	<div><div></div>0.9430</div>	<div><div></div>0.6470</div>

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