



Full wwPDB EM Validation Report ⓘ

Nov 2, 2024 – 10:44 PM EDT

PDB ID : 7MRW
EMDB ID : EMD-23959
Title : Native RhopH complex of the malaria parasite Plasmodium falciparum
Authors : Ho, C.M.; Jih, J.; Lai, M.; Li, X.R.; Goldberg, D.E.; Beck, J.R.; Zhou, Z.H.
Deposited on : 2021-05-10
Resolution : 3.72 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.dev113
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

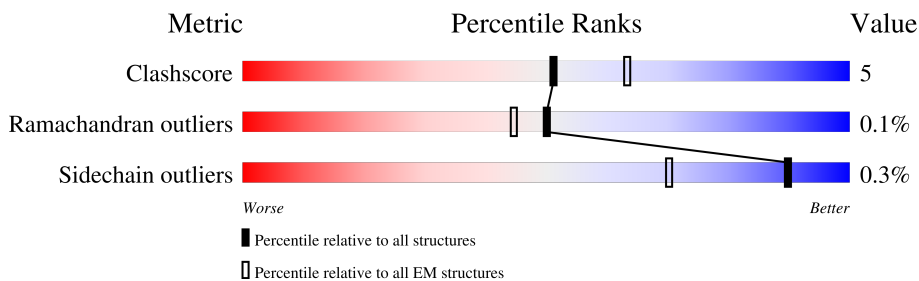
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.72 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	1417	<div> <div>13%</div> <div>71%</div> <div>14%</div> <div>15%</div> </div>
2	B	1378	<div> <div>13%</div> <div>59%</div> <div>12%</div> <div>29%</div> </div>
3	C	897	<div> <div>15%</div> <div>68%</div> <div>11%</div> <div>21%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 24152 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 Cytoadherence linked asexual protein 3.1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1209	Total	C	N	O	S	0	0
			10149	6610	1629	1853	57		

- Molecule 2 is a protein called High molecular weight rhoptry protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	972	Total	C	N	O	S	0	0
			8123	5272	1317	1497	37		

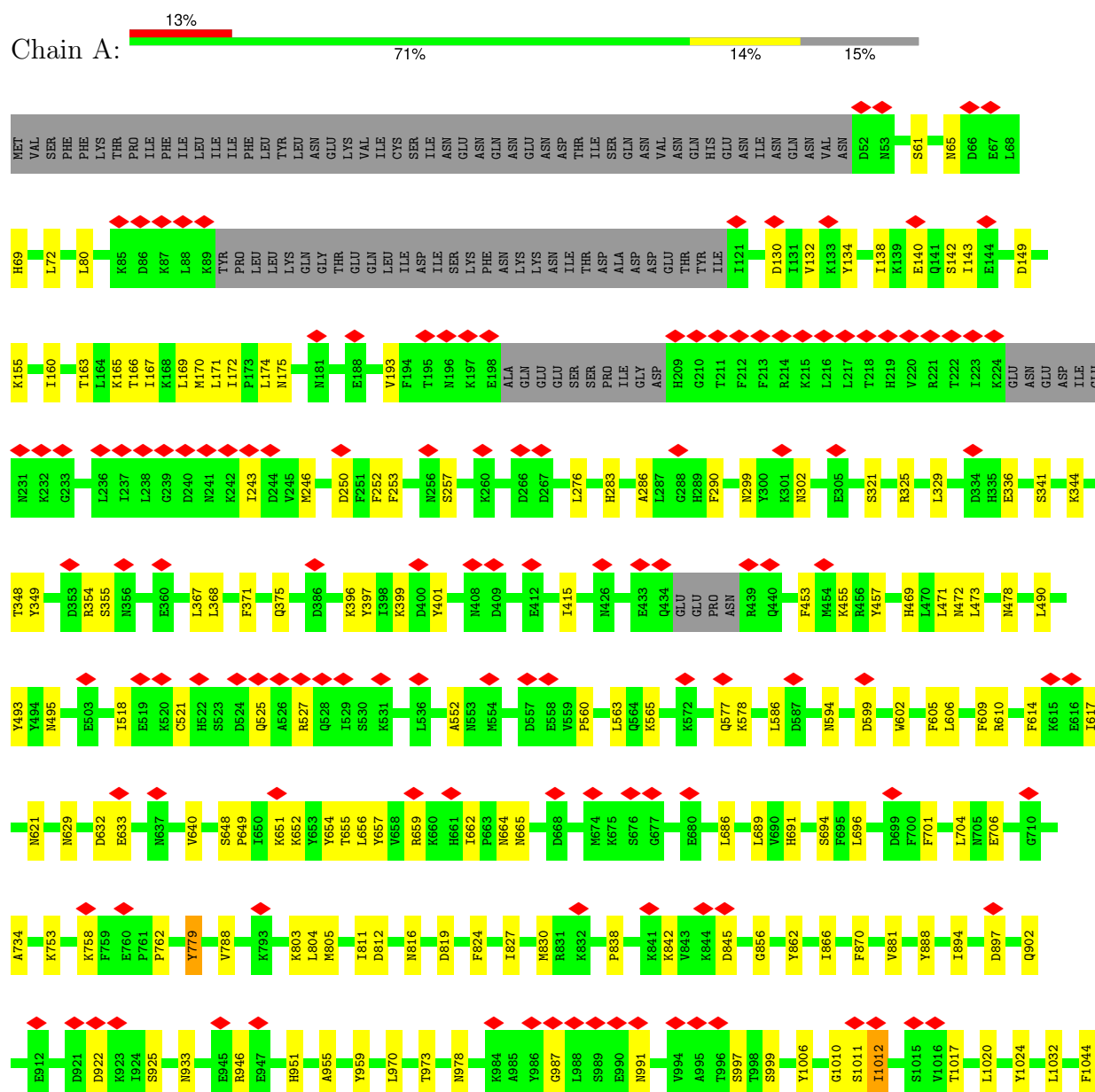
- Molecule 3 is a protein called High molecular weight rhoptry protein 3.

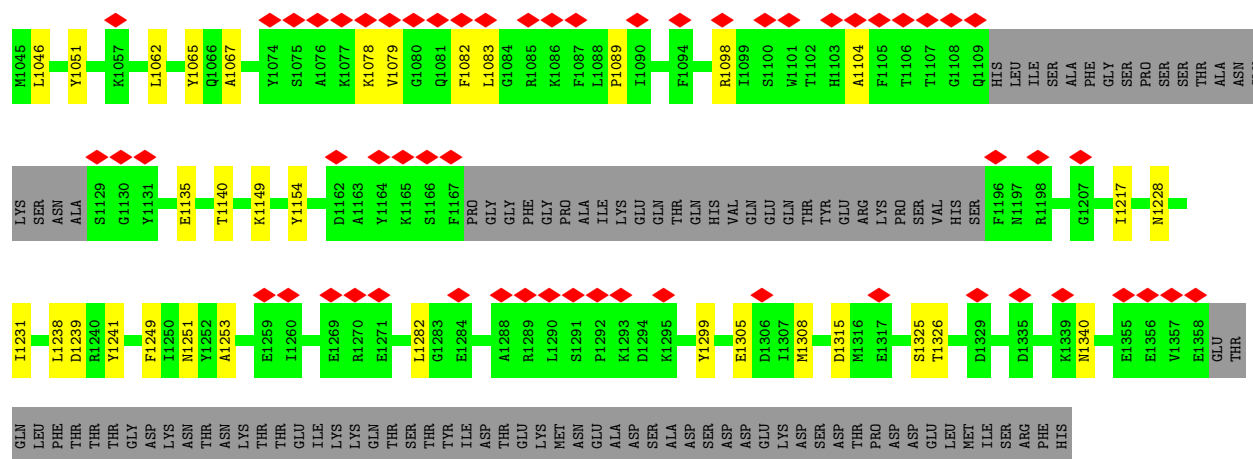
Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	706	Total	C	N	O	S	0	0
			5880	3827	936	1092	25		

3 Residue-property plots

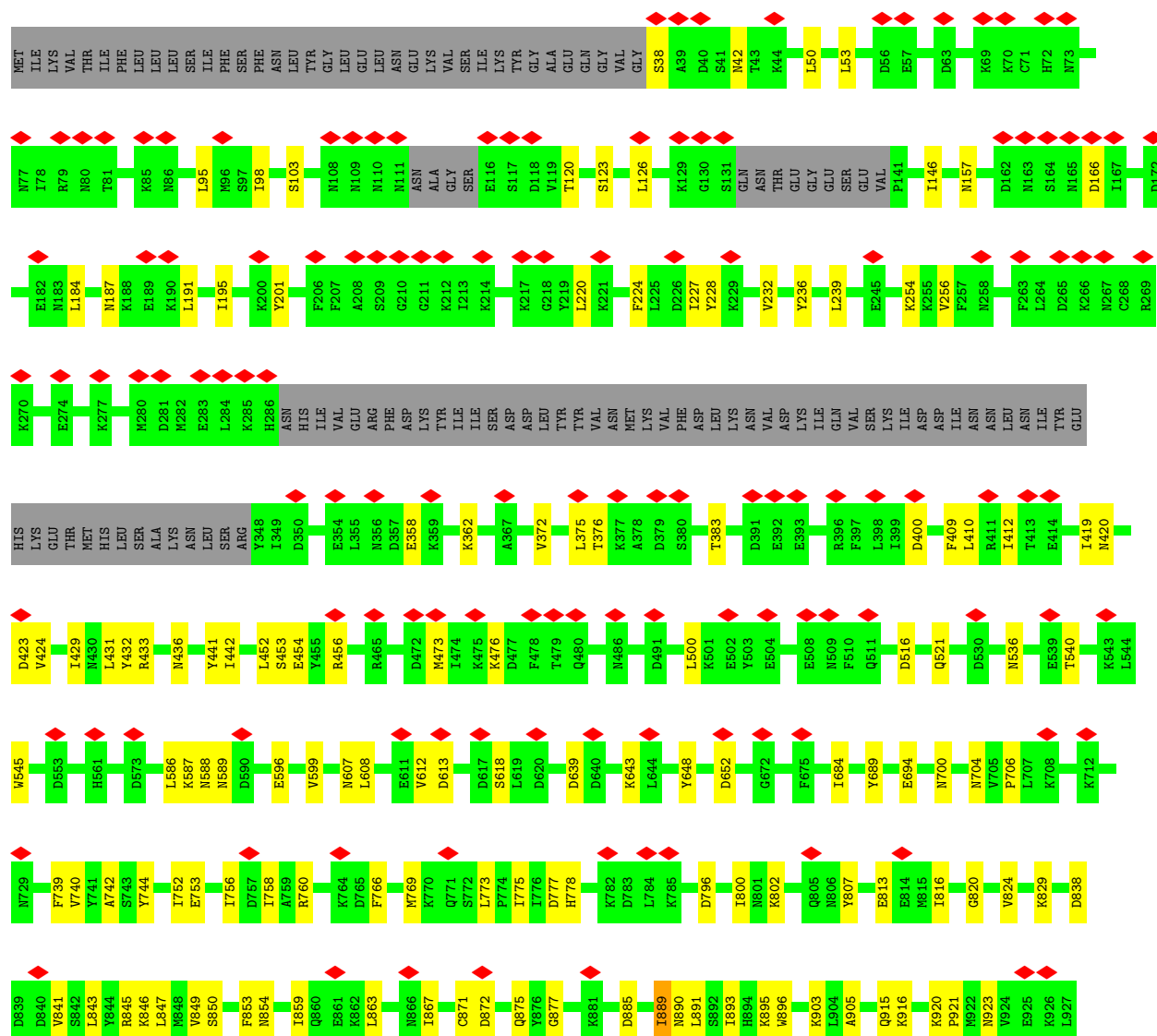
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

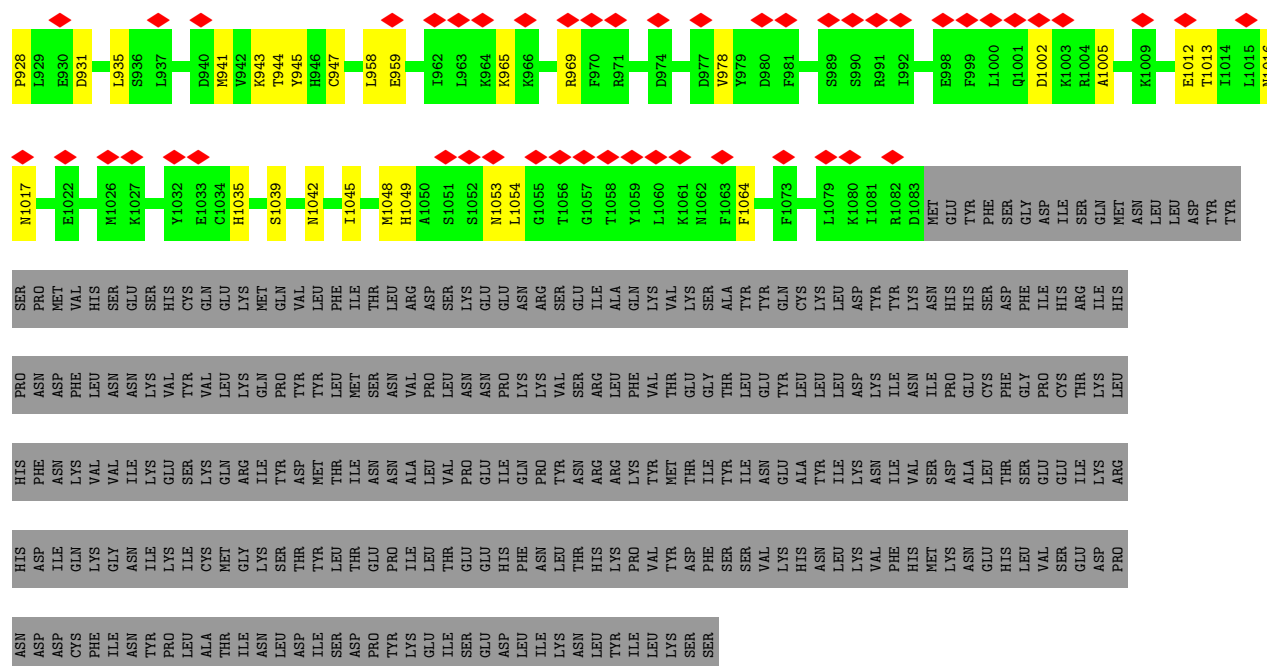
• Molecule 1: Cytoadherence linked asexual protein 3.1



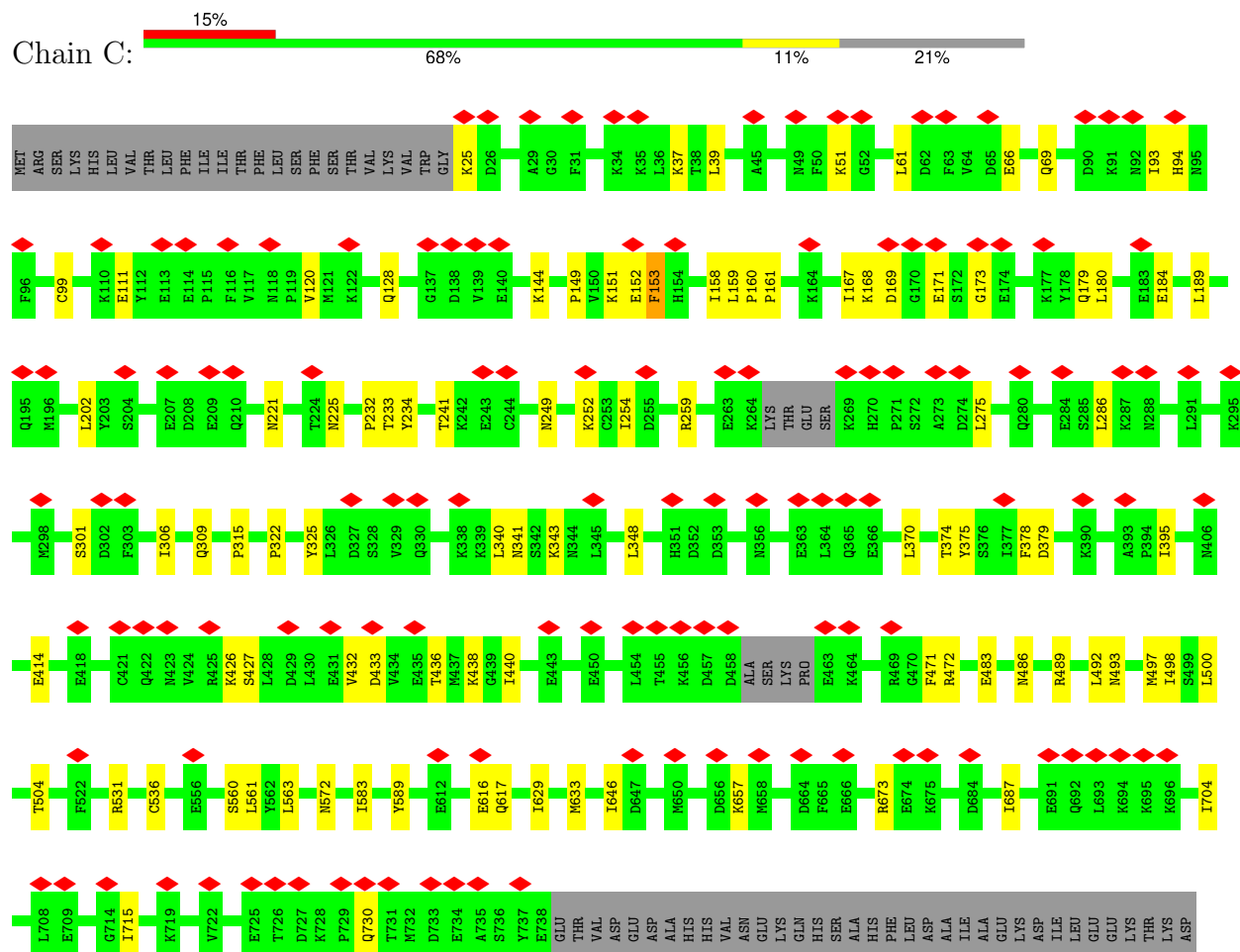


• Molecule 2: High molecular weight rhoptry protein 2





• Molecule 3: High molecular weight rhoptry protein 3



GLU	ASN	GLU	LEU	GLN	LEU	GLU	GLU	ILE	GLU	LEU	TYR	LYS	TYR	MET	GLY	PRO	LEU	LYS	GLU	GLN	SER	LYS	SER	THR	THR	SER	ALA	ALA	SER	THR	THR	SER	ASP	GLU	ILE	SER	GLY	SER	GLY	GLU	PRO	SER	THR	THR	GLU	THR	SER	LYS	THR	THR	GLY	ASN	GLN	GLY	ASP	GLU	ASP	LYS	THR	THR	ASP	ASN	SER	THR	LYS	GLU	MET	GLU
GLU	LEU	GLU	GLU	ALA	GLU	GLY	THR	SER	ASN	LEU	LYS	LYS	GLY	LEU	GLU	PHE	TYR	LYS	SER	SER	LEU	LYS	LEU	ASP	GLN	LEU	SER	ASP	LYS	GLU	LYS	PRO	LYS	SER	LYS	ARG	LYS	LYS	ARG	ASP	SER	SER	SER	ASP	ARG	ILE	LEU	GLU	GLU	SER	LYS	THR	PHE	THR	SER													
GLU	ASN	GLU	LEU																																																																	

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	108872	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.133	Depositor
Minimum map value	-0.086	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.0355	Depositor
Map size (Å)	312.0, 312.0, 312.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.04, 1.04, 1.04	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	0.49	0/10399	0.57	0/14009
2	B	0.43	0/8303	0.53	0/11175
3	C	0.43	0/6025	0.53	0/8130
All	All	0.46	0/24727	0.55	0/33314

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
1	A	0	1
3	C	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1011	SER	Peptide
3	C	286	LEU	Peptide

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	10149	0	10009	116	0
2	B	8123	0	8125	98	0
3	C	5880	0	5830	61	0
All	All	24152	0	23964	257	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 (257) 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:655:THR:HG22	1:A:657:TYR:H	1.55	0.72
3:C:153:PHE:HB2	3:C:348:LEU:HB2	1.71	0.71
2:B:965:LYS:HD2	2:B:969:ARG:HB2	1.76	0.67
3:C:254:ILE:HB	3:C:306:ILE:HG13	1.78	0.65
2:B:410:LEU:HD21	2:B:431:LEU:HD23	1.78	0.65
1:A:1228:ASN:HA	1:A:1231:ILE:HD12	1.78	0.64
3:C:395:ILE:HD11	3:C:657:LYS:HB2	1.80	0.62
2:B:433:ARG:NH2	2:B:586:LEU:O	2.33	0.62
1:A:283:HIS:HD2	1:A:286:ALA:H	1.48	0.61
1:A:1149:LYS:HZ3	3:C:498:ILE:HD11	1.65	0.61
1:A:69:HIS:HB2	1:A:174:LEU:HD22	1.81	0.61
1:A:997:SER:HG	1:A:999:SER:HG	1.47	0.60
2:B:419:ILE:HG23	2:B:423:ASP:HB3	1.83	0.60
1:A:862:TYR:HA	2:B:704:ASN:HD21	1.65	0.60
2:B:166:ASP:OD2	2:B:903:LYS:NZ	2.33	0.60
1:A:140:GLU:HA	1:A:257:SER:HB2	1.84	0.59
1:A:355:SER:HA	1:A:621:ASN:HD21	1.68	0.59
2:B:796:ASP:O	2:B:802:LYS:NZ	2.33	0.59
1:A:881:VAL:HG21	1:A:970:LEU:HD13	1.85	0.59
3:C:432:VAL:O	3:C:489:ARG:NH1	2.36	0.59
2:B:588:ASN:HD21	2:B:694:GLU:HG2	1.68	0.59
1:A:1020:LEU:HD11	1:A:1098:ARG:HG3	1.84	0.58
1:A:706:GLU:OE2	2:B:845:ARG:NH2	2.36	0.58
1:A:649:PRO:HB2	1:A:651:LYS:HE2	1.86	0.58
1:A:838:PRO:O	2:B:441:TYR:OH	2.22	0.58
3:C:315:PRO:HD2	3:C:343:LYS:HE3	1.85	0.57
1:A:987:GLY:HA2	1:A:991:ASN:HB3	1.87	0.57
1:A:321:SER:OG	3:C:572:ASN:ND2	2.38	0.57
1:A:652:LYS:HB2	1:A:894:ILE:HG22	1.86	0.56
2:B:753:GLU:O	2:B:760:ARG:NH2	2.38	0.56
2:B:545:TRP:NE1	2:B:596:GLU:OE2	2.39	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:51:LYS:NZ	3:C:325:TYR:O	2.38	0.56
2:B:959:GLU:OE2	2:B:1053:ASN:ND2	2.38	0.56
2:B:947:CYS:SG	2:B:1035:HIS:N	2.74	0.56
1:A:61:SER:HA	1:A:65:ASN:HB2	1.87	0.55
1:A:704:LEU:O	2:B:689:TYR:OH	2.24	0.55
1:A:689:LEU:H	1:A:870:PHE:HA	1.72	0.55
3:C:69:GLN:NE2	3:C:144:LYS:O	2.39	0.55
3:C:149:PRO:HD2	3:C:340:LEU:HB3	1.88	0.55
3:C:561:LEU:HD12	3:C:633:MET:HG2	1.89	0.55
1:A:518:ILE:HA	1:A:521:CYS:HB3	1.89	0.55
3:C:315:PRO:O	3:C:341:ASN:ND2	2.40	0.55
2:B:126:LEU:HD21	2:B:146:ILE:HD13	1.89	0.54
3:C:120:VAL:HG11	3:C:370:LEU:HD11	1.89	0.54
2:B:228:TYR:HB2	2:B:500:LEU:HD22	1.90	0.54
3:C:438:LYS:HG2	3:C:504:THR:HG23	1.89	0.54
2:B:775:ILE:HD11	2:B:889:ILE:HG13	1.90	0.54
1:A:1006:TYR:O	1:A:1010:GLY:N	2.41	0.54
1:A:838:PRO:HG3	2:B:432:TYR:HE1	1.73	0.53
1:A:577:GLN:HG3	1:A:578:LYS:HG3	1.90	0.53
3:C:379:ASP:OD2	3:C:472:ARG:NH2	2.34	0.53
1:A:453:PHE:HA	1:A:457:TYR:HB2	1.90	0.53
1:A:246:MET:SD	1:A:664:ASN:ND2	2.75	0.53
3:C:168:LYS:NZ	3:C:173:GLY:O	2.40	0.53
1:A:325:ARG:HD3	1:A:368:LEU:HD22	1.91	0.52
1:A:399:LYS:HB2	1:A:478:ASN:HD21	1.74	0.52
1:A:842:LYS:HB3	1:A:845:ASP:HB2	1.91	0.52
2:B:871:CYS:SG	2:B:872:ASP:N	2.83	0.52
2:B:429:ILE:HG13	2:B:589:ASN:HD22	1.74	0.52
1:A:933:ASN:HD22	2:B:587:LYS:HG3	1.75	0.52
2:B:103:SER:OG	2:B:157:ASN:ND2	2.42	0.52
1:A:1140:THR:OG1	1:A:1241:TYR:OH	2.28	0.51
1:A:1340:ASN:ND2	3:C:589:TYR:OH	2.39	0.51
3:C:160:PRO:O	3:C:179:GLN:NE2	2.42	0.51
1:A:686:LEU:O	1:A:951:HIS:NE2	2.44	0.51
2:B:850:SER:O	2:B:854:ASN:ND2	2.44	0.51
1:A:1017:THR:HA	1:A:1020:LEU:HD12	1.92	0.51
1:A:1282:LEU:HD22	3:C:704:ILE:HD13	1.93	0.51
2:B:599:VAL:HG22	2:B:684:ILE:HD11	1.92	0.51
3:C:375:TYR:O	3:C:531:ARG:NH1	2.43	0.51
1:A:138:ILE:O	1:A:142:SER:N	2.41	0.51
2:B:843:LEU:HD13	2:B:896:TRP:CD2	2.46	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1217:ILE:HG21	1:A:1238:LEU:HD13	1.92	0.51
3:C:66:GLU:HG2	3:C:241:THR:HB	1.92	0.51
1:A:734:ALA:HB3	1:A:762:PRO:HD3	1.93	0.50
2:B:187:ASN:HB3	2:B:608:LEU:HD11	1.92	0.50
2:B:372:VAL:HG13	2:B:442:ILE:HD11	1.92	0.50
1:A:701:PHE:HE1	2:B:700:ASN:HD22	1.59	0.50
1:A:525:GLN:OE1	1:A:527:ARG:NH2	2.44	0.50
1:A:640:VAL:HG11	1:A:1299:TYR:HE2	1.76	0.50
2:B:540:THR:HG21	2:B:607:ASN:HB2	1.93	0.50
2:B:739:PHE:HB3	2:B:742:ALA:HB3	1.92	0.50
2:B:890:ASN:HD22	2:B:941:MET:HG2	1.76	0.50
1:A:169:LEU:HD12	1:A:172:ILE:HD12	1.93	0.50
1:A:367:LEU:O	1:A:371:PHE:N	2.44	0.50
2:B:412:ILE:HG21	2:B:419:ILE:HG21	1.93	0.49
3:C:158:ILE:HG13	3:C:159:LEU:HG	1.94	0.49
3:C:616:GLU:OE2	3:C:673:ARG:NH2	2.45	0.49
2:B:859:ILE:HG23	2:B:863:LEU:HD13	1.93	0.49
1:A:160:ILE:O	1:A:163:THR:OG1	2.28	0.49
1:A:1079:VAL:HA	1:A:1083:LEU:HD12	1.93	0.49
3:C:259:ARG:HH12	3:C:301:SER:HA	1.78	0.49
1:A:803:LYS:NZ	2:B:694:GLU:OE1	2.40	0.49
2:B:1012:GLU:O	2:B:1016:ASN:ND2	2.46	0.49
2:B:201:TYR:HB3	2:B:232:VAL:HG23	1.93	0.49
2:B:807:TYR:OH	2:B:813:GLU:OE2	2.28	0.49
3:C:560:SER:HB2	3:C:563:LEU:HB2	1.95	0.49
2:B:195:ILE:HG23	2:B:536:ASN:HD21	1.77	0.48
3:C:492:LEU:HG	3:C:629:ILE:HD11	1.93	0.48
1:A:246:MET:HB3	1:A:662:ILE:HD13	1.95	0.48
1:A:149:ASP:OD1	1:A:654:TYR:OH	2.31	0.48
1:A:1326:THR:OG1	3:C:498:ILE:O	2.26	0.48
3:C:151:LYS:HG2	3:C:233:THR:HG22	1.96	0.48
1:A:299:ASN:HB3	1:A:302:ASN:HB2	1.96	0.48
1:A:143:ILE:HG12	1:A:253:PHE:CD2	2.49	0.48
1:A:472:ASN:O	1:A:493:TYR:OH	2.32	0.48
1:A:632:ASP:OD1	1:A:632:ASP:N	2.46	0.48
3:C:25:LYS:N	3:C:111:GLU:O	2.46	0.48
1:A:1089:PRO:HG3	1:A:1154:TYR:CZ	2.49	0.48
2:B:453:SER:OG	2:B:454:GLU:N	2.46	0.48
2:B:889:ILE:HG21	2:B:945:TYR:HD1	1.78	0.48
3:C:583:ILE:HD11	3:C:646:ILE:HG23	1.94	0.48
1:A:779:TYR:CD2	1:A:811:ILE:HG12	2.49	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:LEU:HD11	1:A:290:PHE:HB3	1.96	0.47
1:A:354:ARG:O	1:A:621:ASN:ND2	2.47	0.47
1:A:659:ARG:HD2	1:A:665:ASN:HD22	1.79	0.47
2:B:358:GLU:HG3	2:B:362:LYS:HE2	1.96	0.47
2:B:769:MET:HA	2:B:843:LEU:HD23	1.95	0.47
1:A:599:ASP:HB3	1:A:602:TRP:CD1	2.48	0.47
2:B:412:ILE:HG13	2:B:424:VAL:HG22	1.96	0.47
3:C:39:LEU:HD23	3:C:99:CYS:HB2	1.96	0.47
1:A:172:ILE:HD13	1:A:193:VAL:HG21	1.96	0.47
2:B:95:LEU:HD23	2:B:98:ILE:HD12	1.97	0.47
2:B:846:LYS:NZ	2:B:895:LYS:O	2.38	0.47
3:C:471:PHE:HZ	3:C:536:CYS:HB2	1.79	0.47
3:C:152:GLU:OE2	3:C:309:GLN:NE2	2.42	0.47
1:A:140:GLU:HG2	1:A:257:SER:HB2	1.97	0.47
2:B:1039:SER:OG	2:B:1042:ASN:ND2	2.48	0.47
1:A:495:ASN:ND2	1:A:648:SER:OG	2.48	0.47
2:B:1045:ILE:HA	2:B:1048:MET:HB2	1.97	0.47
1:A:336:GLU:H	1:A:354:ARG:HH12	1.61	0.46
1:A:341:SER:HB2	1:A:348:THR:HG21	1.96	0.46
1:A:1249:PHE:O	1:A:1253:ALA:N	2.48	0.46
1:A:1012:ILE:H	1:A:1012:ILE:HG13	1.55	0.46
3:C:189:LEU:HD23	3:C:715:ILE:HD11	1.97	0.46
3:C:249:ASN:HB2	3:C:252:LYS:HG2	1.98	0.46
3:C:497:MET:HA	3:C:500:LEU:HG	1.98	0.46
1:A:375:GLN:HA	1:A:473:LEU:HD11	1.97	0.46
1:A:696:LEU:HD13	1:A:866:ILE:HG23	1.97	0.46
2:B:452:LEU:HD13	2:B:456:ARG:HD3	1.98	0.46
2:B:816:ILE:O	2:B:820:GLY:N	2.44	0.46
2:B:409:PHE:HB2	2:B:424:VAL:HG12	1.98	0.45
2:B:838:ASP:HB3	2:B:841:VAL:HG12	1.98	0.45
1:A:132:VAL:HG22	1:A:165:LYS:HD2	1.97	0.45
1:A:252:PHE:O	1:A:602:TRP:NE1	2.38	0.45
2:B:652:ASP:O	2:B:829:LYS:NZ	2.49	0.45
1:A:472:ASN:HD22	1:A:490:LEU:HD11	1.80	0.45
2:B:236:TYR:HB3	2:B:239:LEU:HB2	1.98	0.45
2:B:224:PHE:HD1	2:B:227:ILE:HD11	1.81	0.45
2:B:893:ILE:HG13	2:B:944:THR:HG21	1.99	0.45
3:C:169:ASP:HB3	3:C:171:GLU:HG2	1.99	0.45
1:A:552:ALA:HB1	1:A:563:LEU:HG	1.97	0.45
2:B:700:ASN:ND2	2:B:706:PRO:O	2.48	0.45
2:B:778:HIS:NE2	2:B:885:ASP:OD1	2.50	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:483:GLU:OE2	3:C:504:THR:N	2.49	0.45
1:A:490:LEU:HD12	1:A:493:TYR:HE2	1.82	0.45
1:A:344:LYS:NZ	1:A:633:GLU:OE1	2.45	0.45
2:B:254:LYS:NZ	2:B:516:ASP:OD2	2.49	0.45
2:B:773:LEU:O	2:B:896:TRP:NE1	2.49	0.45
1:A:72:LEU:HD22	1:A:167:ILE:HD12	1.99	0.45
1:A:973:THR:HG21	1:A:1051:TYR:HE1	1.81	0.45
3:C:486:ASN:ND2	3:C:504:THR:OG1	2.46	0.45
1:A:691:HIS:O	1:A:694:SER:OG	2.35	0.44
1:A:606:LEU:O	1:A:610:ARG:N	2.40	0.44
2:B:777:ASP:OD2	2:B:1049:HIS:NE2	2.50	0.44
2:B:38:SER:O	2:B:42:ASN:N	2.48	0.44
2:B:773:LEU:H	2:B:896:TRP:HE1	1.64	0.44
2:B:875:GLN:HE22	2:B:915:GLN:HB3	1.83	0.44
1:A:605:PHE:O	1:A:609:PHE:N	2.51	0.44
3:C:161:PRO:HG2	3:C:180:LEU:HD12	1.98	0.44
2:B:800:ILE:HB	2:B:813:GLU:HG3	1.99	0.44
2:B:875:GLN:HA	2:B:921:PRO:HB3	2.00	0.44
2:B:935:LEU:HD11	2:B:943:LYS:HA	1.99	0.44
3:C:61:LEU:H	3:C:128:GLN:HE22	1.65	0.44
3:C:374:THR:O	3:C:378:PHE:N	2.46	0.44
3:C:322:PRO:HA	3:C:325:TYR:HD2	1.82	0.44
1:A:396:LYS:NZ	3:C:414:GLU:OE2	2.50	0.44
1:A:888:TYR:OH	1:A:897:ASP:OD1	2.29	0.44
1:A:1044:PHE:HB2	1:A:1067:ALA:HB2	2.00	0.44
3:C:61:LEU:HD22	3:C:69:GLN:HB3	1.99	0.44
1:A:614:PHE:HB2	1:A:617:ILE:HB	2.00	0.44
1:A:788:VAL:HG13	1:A:804:LEU:HD22	2.00	0.44
2:B:50:LEU:HD23	2:B:53:LEU:HD12	2.00	0.44
2:B:376:THR:O	2:B:383:THR:OG1	2.35	0.44
2:B:824:VAL:HG12	2:B:849:VAL:HG22	2.00	0.43
2:B:473:MET:HA	2:B:476:LYS:HE2	2.00	0.43
2:B:648:TYR:HD1	2:B:853:PHE:HE2	1.67	0.43
2:B:958:LEU:HD13	2:B:978:VAL:HG11	2.00	0.43
2:B:1013:THR:O	2:B:1017:ASN:ND2	2.51	0.43
1:A:155:LYS:NZ	1:A:594:ASN:O	2.50	0.43
1:A:349:TYR:CE1	1:A:354:ARG:HG3	2.53	0.43
1:A:824:PHE:HZ	1:A:1046:LEU:HD22	1.84	0.43
2:B:120:THR:O	2:B:123:SER:OG	2.31	0.43
2:B:375:LEU:HD12	2:B:442:ILE:HD12	1.99	0.43
2:B:752:ILE:HG12	2:B:756:ILE:HD11	1.99	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:920:LYS:HD2	2:B:921:PRO:HD2	2.00	0.43
1:A:329:LEU:HD22	1:A:368:LEU:HD12	2.00	0.43
1:A:1104:ALA:HB3	1:A:1135:GLU:HB3	2.01	0.43
2:B:400:ASP:OD2	2:B:436:ASN:ND2	2.49	0.43
2:B:184:LEU:HD21	2:B:612:VAL:HG13	2.01	0.43
2:B:256:VAL:HG21	2:B:442:ILE:HG21	2.00	0.43
2:B:928:PRO:HD2	2:B:931:ASP:HB2	2.00	0.43
1:A:250:ASP:O	1:A:565:LYS:NZ	2.45	0.43
1:A:955:ALA:HB1	1:A:959:TYR:HB2	2.00	0.43
2:B:545:TRP:CD1	2:B:740:VAL:HB	2.53	0.43
1:A:469:HIS:O	1:A:473:LEU:N	2.52	0.43
1:A:656:LEU:HD13	1:A:902:GLN:HG2	2.01	0.43
1:A:819:ASP:OD1	1:A:819:ASP:N	2.52	0.43
2:B:613:ASP:OD1	2:B:613:ASP:N	2.52	0.43
1:A:171:LEU:O	1:A:175:ASN:ND2	2.52	0.43
1:A:415:ILE:HD12	1:A:471:LEU:HD22	2.00	0.42
1:A:830:MET:HB3	1:A:856:GLY:HA3	2.01	0.42
1:A:397:TYR:O	1:A:401:TYR:N	2.45	0.42
1:A:1305:GLU:HA	1:A:1308:MET:HB2	2.01	0.42
2:B:847:LEU:HD23	2:B:847:LEU:HA	1.83	0.42
3:C:184:GLU:HB3	3:C:202:LEU:HD11	2.00	0.42
1:A:1251:ASN:ND2	3:C:493:ASN:OD1	2.52	0.42
1:A:166:THR:O	1:A:170:MET:N	2.46	0.42
1:A:827:ILE:HG12	1:A:1032:LEU:HD21	2.01	0.42
3:C:234:TYR:CE2	3:C:275:LEU:HB2	2.54	0.42
1:A:812:ASP:O	1:A:816:ASN:ND2	2.53	0.42
1:A:1062:LEU:HA	1:A:1065:TYR:HD2	1.85	0.42
1:A:1325:SER:HB3	3:C:500:LEU:HD12	2.02	0.42
1:A:978:ASN:HD22	1:A:1239:ASP:HA	1.85	0.42
1:A:1024:TYR:OH	1:A:1315:ASP:OD2	2.25	0.42
1:A:130:ASP:O	1:A:134:TYR:N	2.53	0.42
2:B:877:GLY:HA3	2:B:923:ASN:HA	2.02	0.42
3:C:440:ILE:HG21	3:C:687:ILE:HG12	2.00	0.42
1:A:1078:LYS:HG2	1:A:1082:PHE:HE1	1.85	0.42
2:B:191:LEU:HD13	2:B:608:LEU:HD22	2.01	0.42
2:B:639:ASP:OD1	2:B:639:ASP:N	2.50	0.42
3:C:563:LEU:HD23	3:C:563:LEU:HA	1.82	0.42
2:B:775:ILE:HD11	2:B:889:ILE:HA	2.00	0.41
3:C:433:ASP:O	3:C:436:THR:OG1	2.25	0.41
3:C:93:ILE:HG22	3:C:94:HIS:H	1.85	0.41
1:A:862:TYR:HA	2:B:704:ASN:ND2	2.33	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:872:ASP:OD2	2:B:916:LYS:NZ	2.50	0.41
3:C:161:PRO:HA	3:C:232:PRO:HD3	2.01	0.41
1:A:922:ASP:N	1:A:925:SER:OG	2.42	0.41
2:B:618:SER:HA	2:B:643:LYS:HG2	2.03	0.41
2:B:220:LEU:O	2:B:521:GLN:NE2	2.54	0.41
2:B:758:ILE:HD11	2:B:766:PHE:CD2	2.56	0.41
3:C:426:LYS:HA	3:C:617:GLN:HE22	1.84	0.41
1:A:753:LYS:O	1:A:758:LYS:N	2.51	0.41
1:A:243:ILE:HD13	1:A:560:PRO:HD3	2.03	0.41
1:A:805:MET:SD	1:A:946:ARG:NH1	2.93	0.41
1:A:1078:LYS:HD2	3:C:730:GLN:HE21	1.85	0.41
2:B:867:ILE:HD11	2:B:905:ALA:HB1	2.03	0.41
3:C:427:SER:H	3:C:617:GLN:HE22	1.69	0.41
3:C:161:PRO:O	3:C:179:GLN:NE2	2.53	0.41
3:C:167:ILE:HG12	3:C:225:ASN:HD22	1.86	0.40
2:B:744:TYR:HD1	2:B:744:TYR:HA	1.77	0.40
1:A:80:LEU:HD23	1:A:80:LEU:HA	1.92	0.40
3:C:153:PHE:HB3	3:C:158:ILE:HG22	2.02	0.40
3:C:221:ASN:HB3	3:C:225:ASN:HD22	1.86	0.40
2:B:1002:ASP:HB3	2:B:1005:ALA:HB3	2.03	0.40
2:B:1054:LEU:HB3	2:B:1064:PHE:HE2	1.87	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	1195/1417 (84%)	1146 (96%)	47 (4%)	2 (0%)	44	72
2	B	964/1378 (70%)	934 (97%)	30 (3%)	0	100	100
3	C	700/897 (78%)	668 (95%)	32 (5%)	0	100	100
All	All	2859/3692 (77%)	2748 (96%)	109 (4%)	2 (0%)	50	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	586	LEU
1	A	1012	ILE

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	1129/1321 (86%)	1126 (100%)	3 (0%)	91	94
2	B	915/1301 (70%)	912 (100%)	3 (0%)	91	94
3	C	672/847 (79%)	670 (100%)	2 (0%)	91	94
All	All	2716/3469 (78%)	2708 (100%)	8 (0%)	90	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	455	LYS
1	A	629	ASN
1	A	779	TYR
2	B	420	ASN
2	B	889	ILE
2	B	891	LEU
3	C	37	LYS
3	C	153	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	175	ASN
1	A	279	HIS
1	A	283	HIS
1	A	472	ASN
1	A	478	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	621	ASN
1	A	730	ASN
1	A	816	ASN
1	A	952	ASN
1	A	968	GLN
1	A	978	ASN
1	A	1251	ASN
1	A	1301	HIS
1	A	1340	ASN
2	B	42	ASN
2	B	152	ASN
2	B	157	ASN
2	B	286	HIS
2	B	420	ASN
2	B	450	ASN
2	B	588	ASN
2	B	635	HIS
2	B	666	ASN
2	B	875	GLN
2	B	890	ASN
2	B	1017	ASN
2	B	1030	ASN
2	B	1042	ASN
3	C	225	ASN
3	C	351	HIS
3	C	468	HIS
3	C	486	ASN
3	C	572	ASN
3	C	617	GLN
3	C	637	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

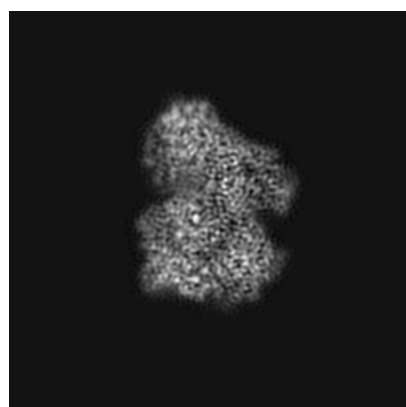
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23959. These allow visual inspection of the internal detail of the map and identification of artifacts.

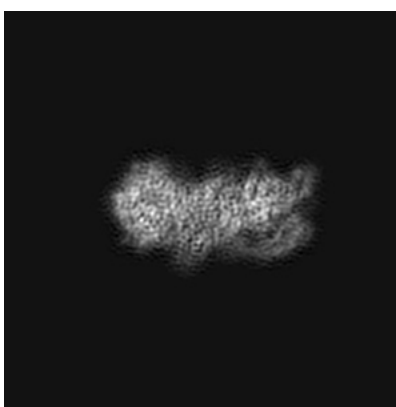
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

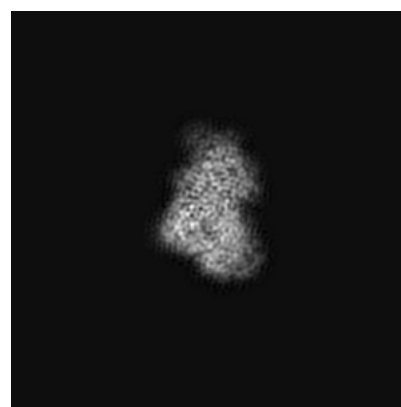
6.1.1 Primary map



X



Y

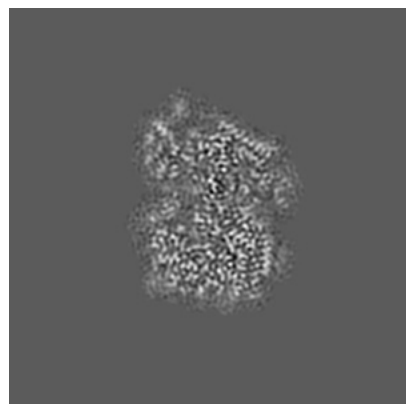


Z

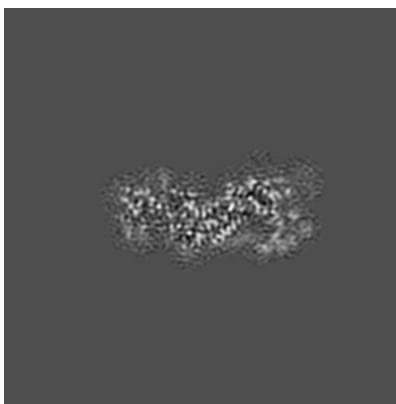
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

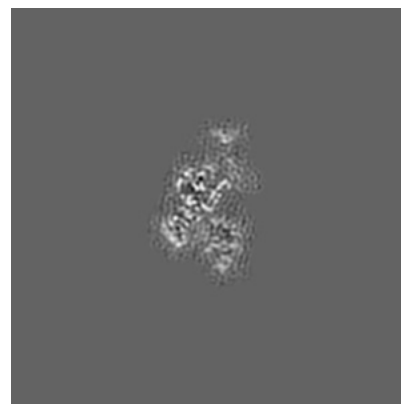
6.2.1 Primary map



X Index: 150



Y Index: 150

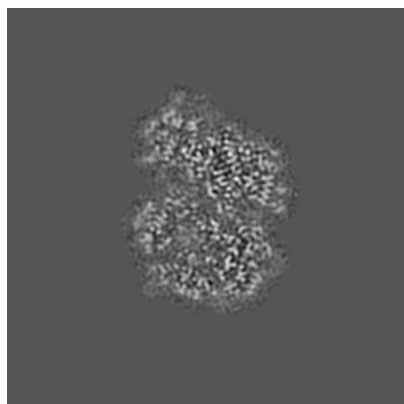


Z Index: 150

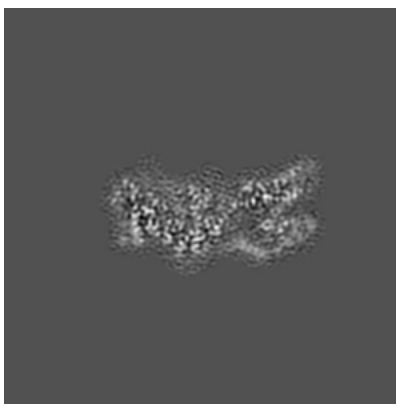
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

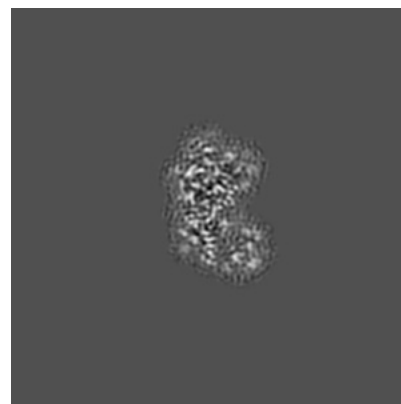
6.3.1 Primary map



X Index: 156



Y Index: 140

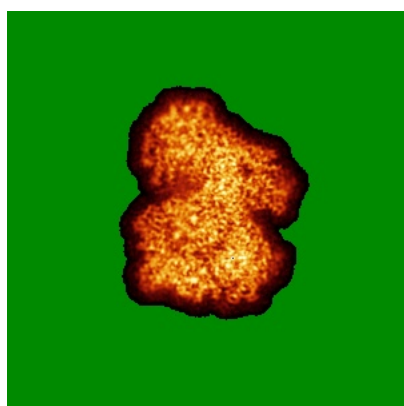


Z Index: 110

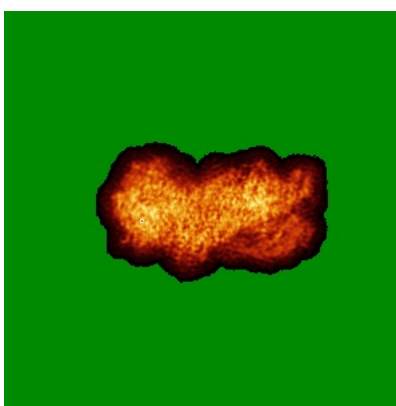
The images above show the largest variance slices of the map in three orthogonal directions.

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

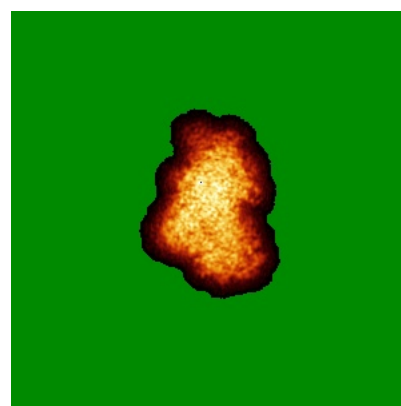
6.4.1 Primary map



X



Y

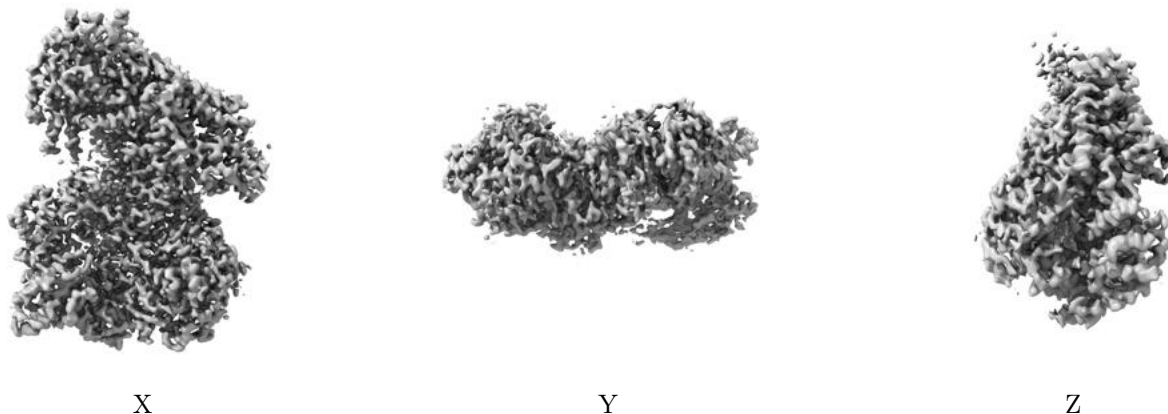


Z

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 0.0355. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

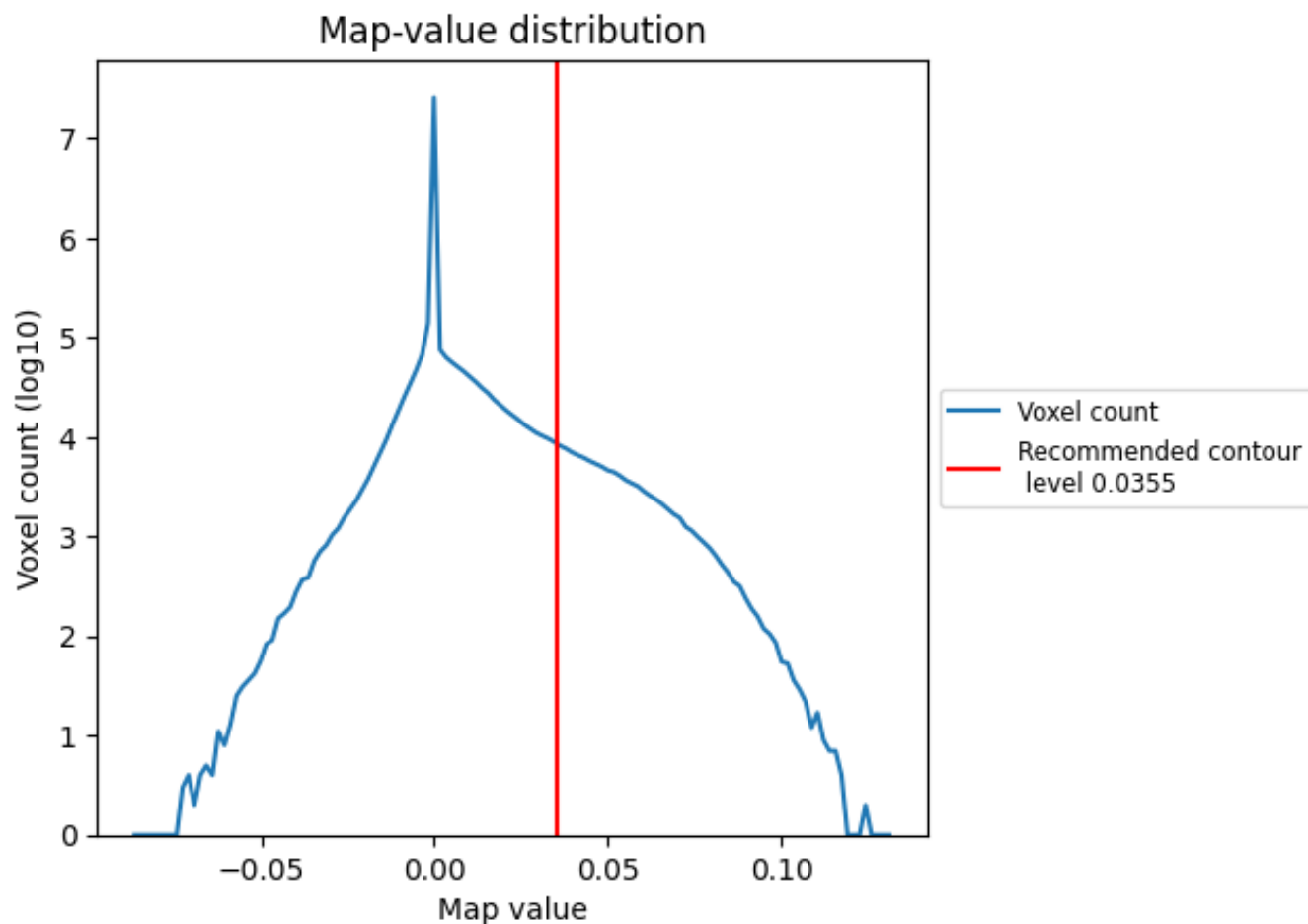
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

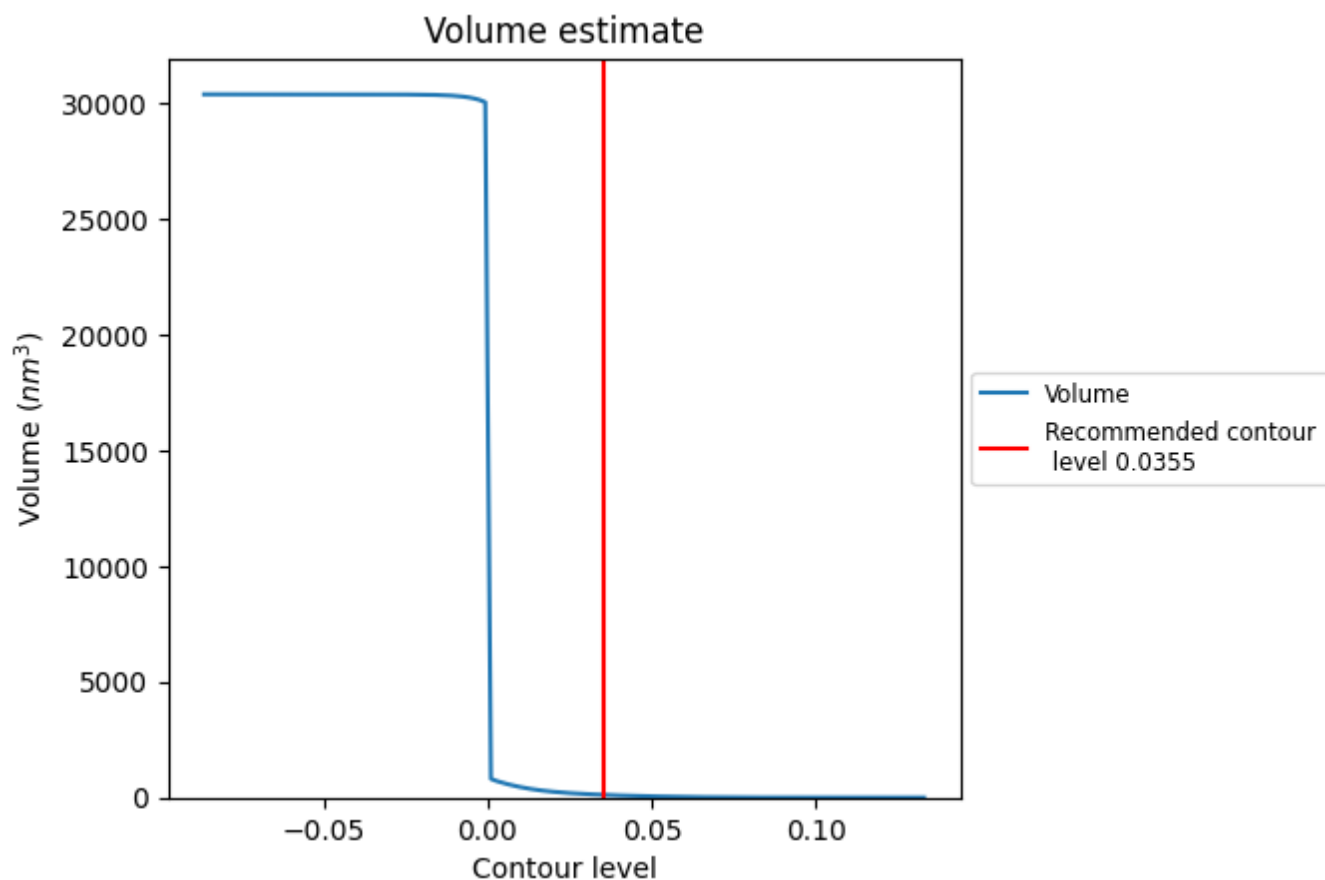
This section contains the results of statistical analysis of the map.

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.

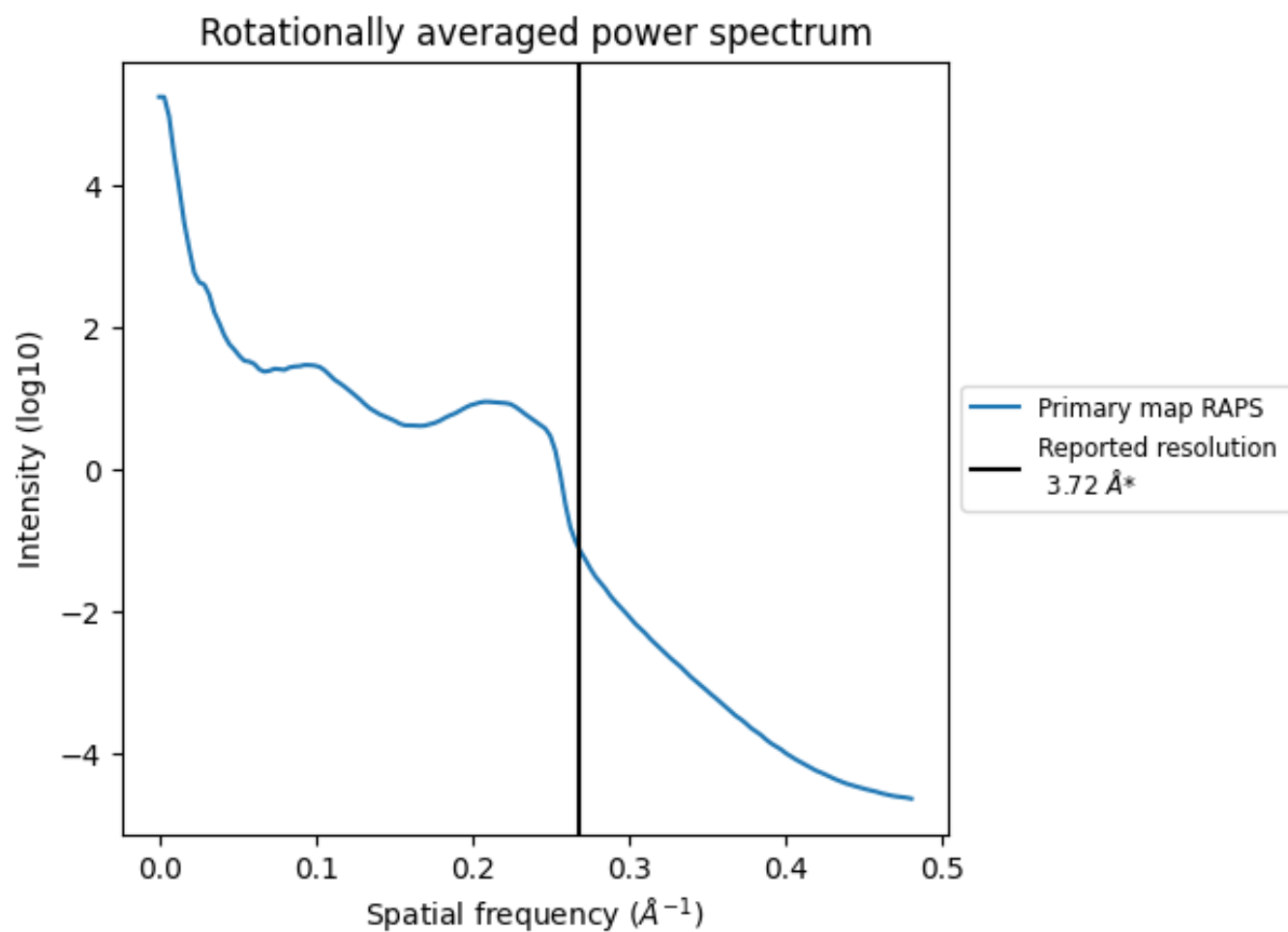
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 115 nm³; this corresponds to an approximate mass of 104 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 0.269 Å⁻¹

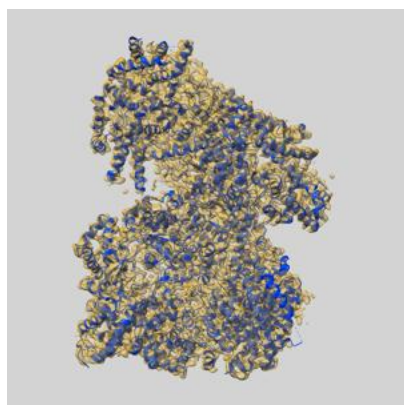
8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

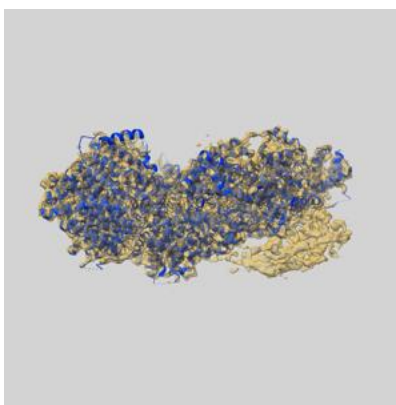
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23959 and PDB model 7MRW. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

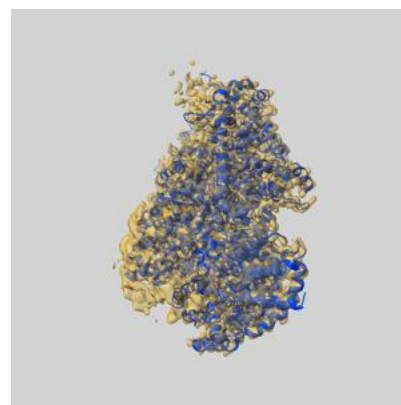
9.1 Map-model overlay [i](#)



X



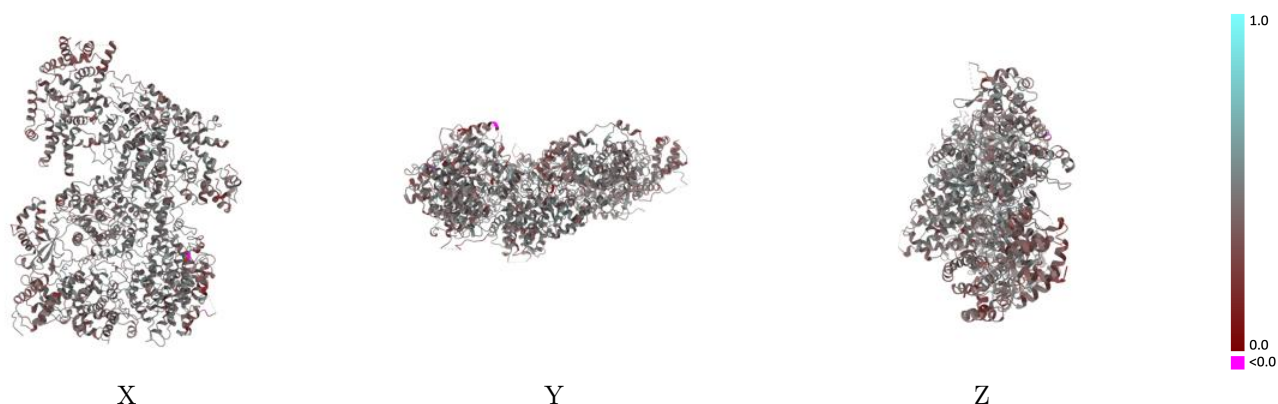
Y



Z

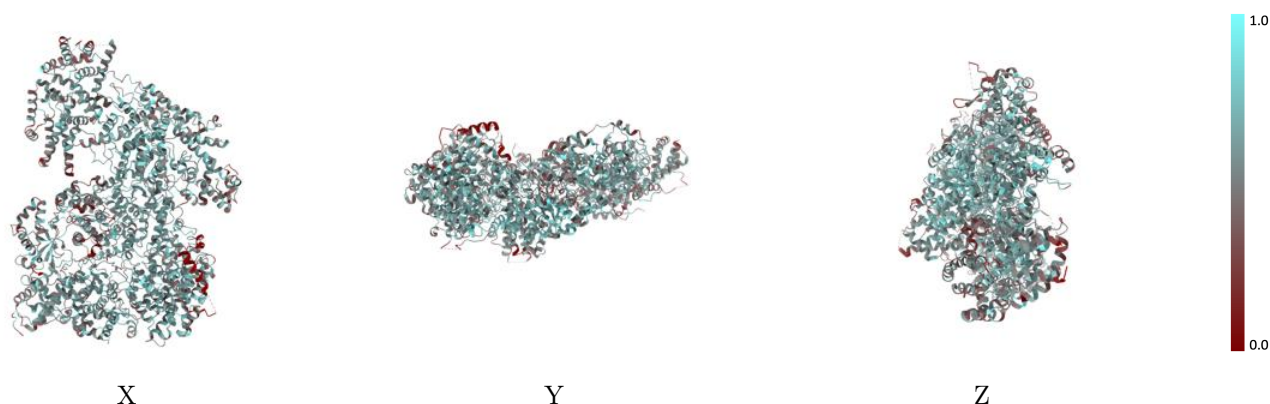
The images above show the 3D surface view of the map at the recommended contour level 0.0355 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



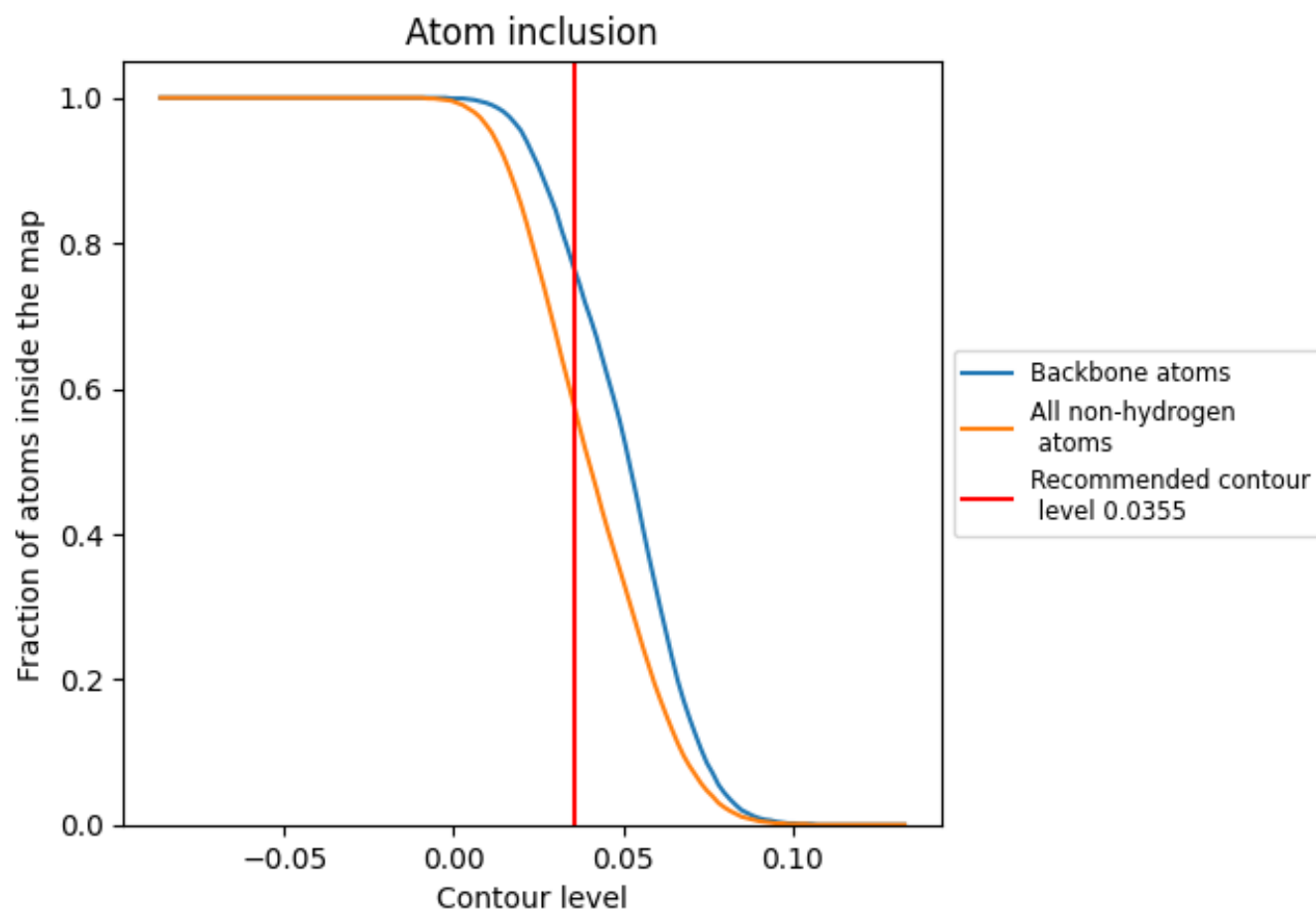
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0355).

9.4 Atom inclusion [i](#)



At the recommended contour level, 77% of all backbone atoms, 58% 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 (0.0355) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.5770	<div></div> 0.4320
A	<div></div> 0.5930	<div></div> 0.4430
B	<div></div> 0.5630	<div></div> 0.4240
C	<div></div> 0.5680	<div></div> 0.4220

