



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 07:08 AM UTC

PDB ID : 9M0Y / pdb_00009m0y
EMDB ID : EMD-63565
Title : Local refinement of stacked like DDB1-DDA1-DET1-Ube2e2-COP1 complex (layer 2)
Authors : Su, M.-Y.
Deposited on : 2025-02-25
Resolution : 4.25 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

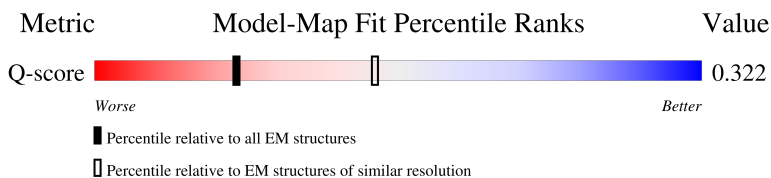
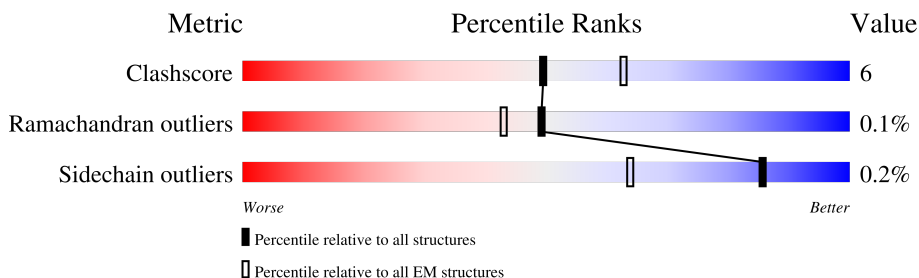
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 4.25 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.



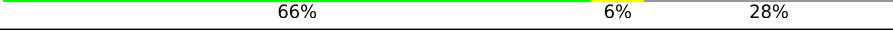
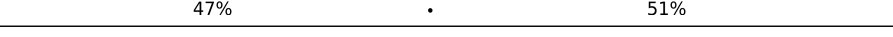
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	4654 (3.75 - 4.75)

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	731	
1	B	731	
1	C	731	
1	D	731	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	E	731	
1	F	731	
1	G	731	
1	J	731	
2	Q	201	
2	T	201	
3	R	550	
3	U	550	
4	S	1140	
4	V	1140	
5	j	102	
5	n	102	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 36013 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 E3 ubiquitin-protein ligase COP1.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	434	Total	C	N	O	0	0
			2213	1334	443	436		
1	B	434	Total	C	N	O	0	0
			2182	1313	435	434		
1	C	532	Total	C	N	O	0	0
			2715	1640	539	536		
1	D	503	Total	C	N	O	0	0
			2519	1513	503	503		
1	E	527	Total	C	N	O	0	0
			2704	1632	540	532		
1	F	508	Total	C	N	O	0	0
			2544	1528	508	508		
1	G	435	Total	C	N	O	0	0
			2173	1303	435	435		
1	J	435	Total	C	N	O	0	0
			2173	1303	435	435		

- Molecule 2 is a protein called Ubiquitin-conjugating enzyme E2 E2.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	Q	147	Total	C	N	O	0	0
			786	488	149	149		
2	T	147	Total	C	N	O	0	0
			811	506	155	150		

- Molecule 3 is a protein called DET1 homolog.

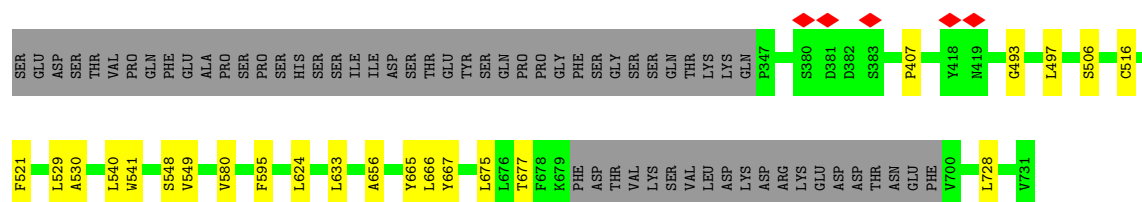
Mol	Chain	Residues	Atoms					AltConf	Trace
3	R	522	Total	C	N	O		0	0
			2976	1854	576	546			
3	U	540	Total	C	N	O	S	0	0
			3038	1881	588	568	1		

- Molecule 4 is a protein called DNA damage-binding protein 1.

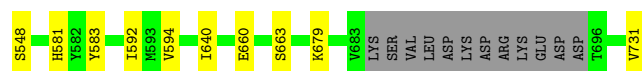
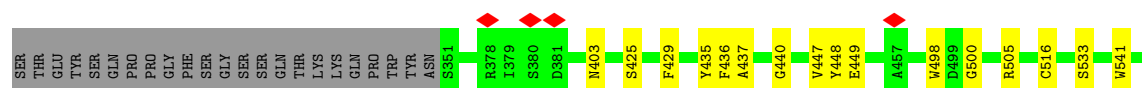
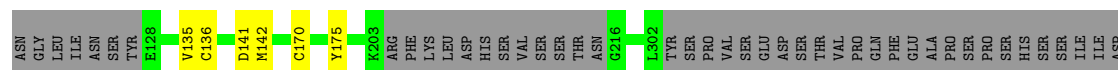
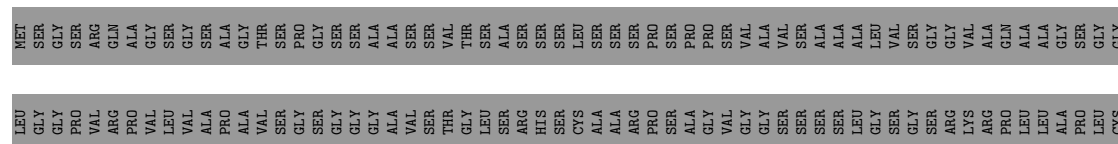
Mol	Chain	Residues	Atoms					AltConf	Trace
4	S	823	Total	C	N	O	S	0	0
			4313	2621	854	835	3		
4	V	823	Total	C	N	O		0	0
			4234	2555	844	835			

- Molecule 5 is a protein called DET1- and DDB1-associated protein 1.

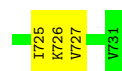
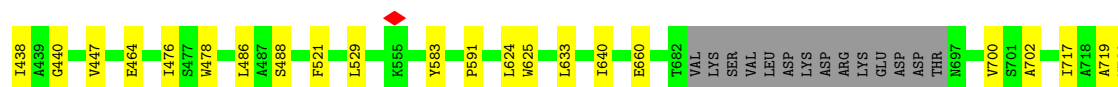
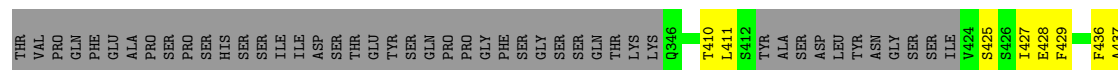
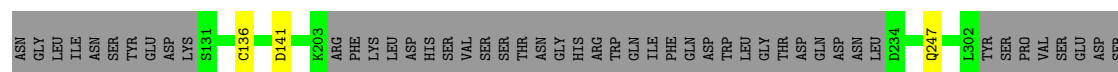
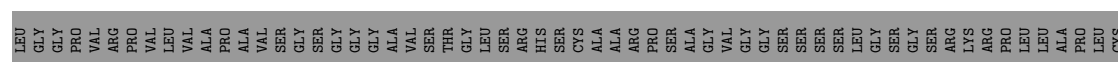
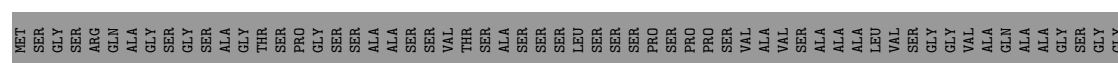
Mol	Chain	Residues	Atoms				AltConf	Trace
5	j	50	Total	C	N	O	0	0
			299	186	58	55		
5	n	49	Total	C	N	O	0	0
			333	214	59	60		



• Molecule 1: E3 ubiquitin-protein ligase COP1

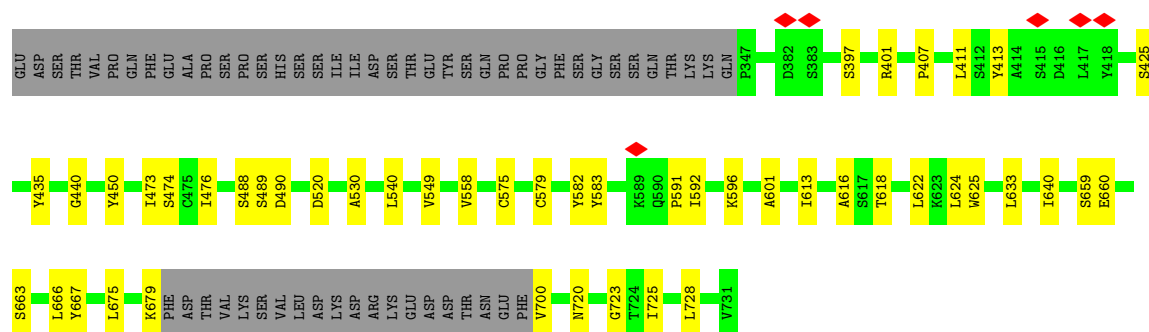


• Molecule 1: E3 ubiquitin-protein ligase COP1



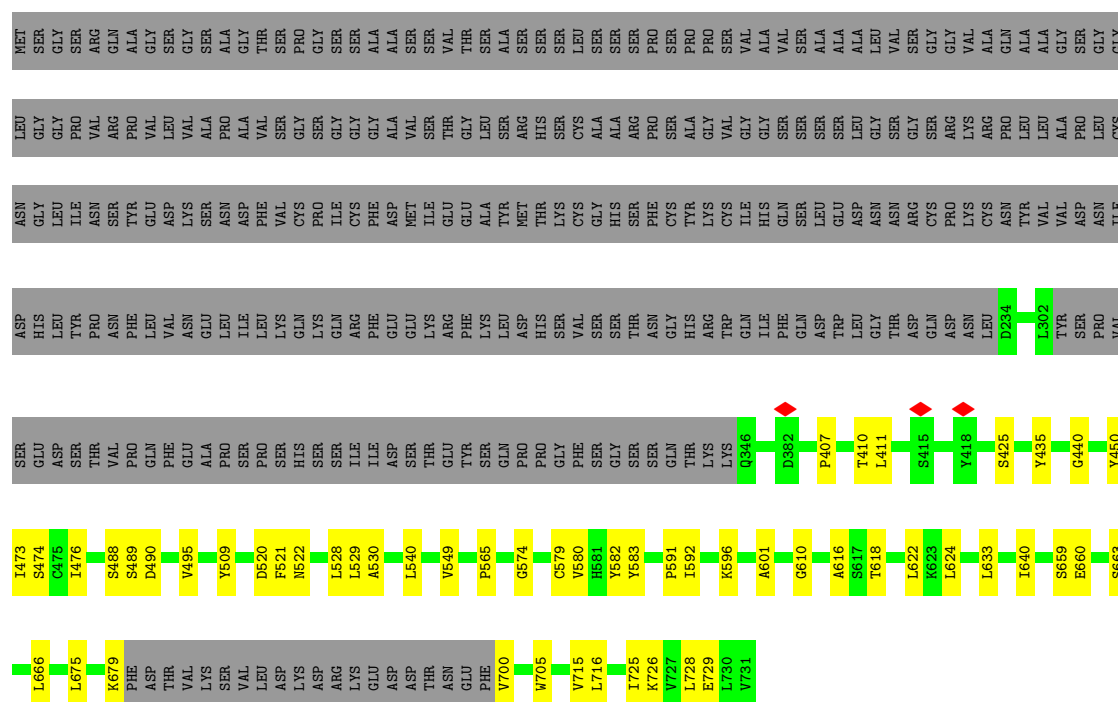
• Molecule 1: E3 ubiquitin-protein ligase COP1





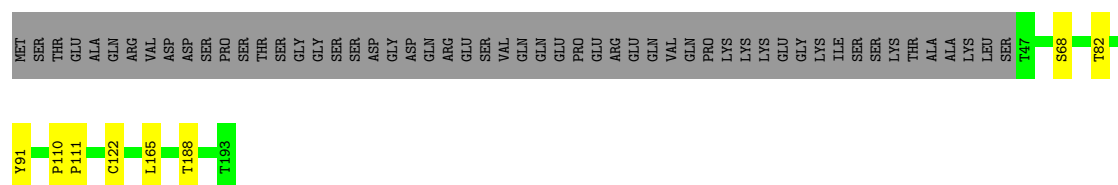
• Molecule 1: E3 ubiquitin-protein ligase COP1

Chain J: 52% 7% 40%



• Molecule 2: Ubiquitin-conjugating enzyme E2 E2

Chain Q: 69% 27%

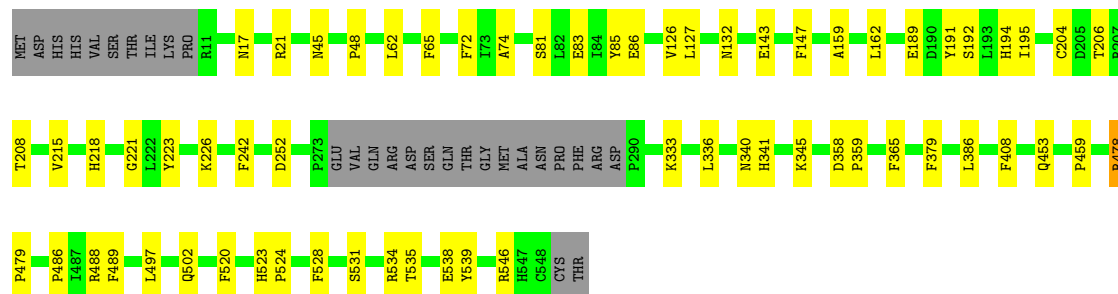
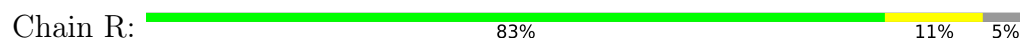


• Molecule 2: Ubiquitin-conjugating enzyme E2 E2

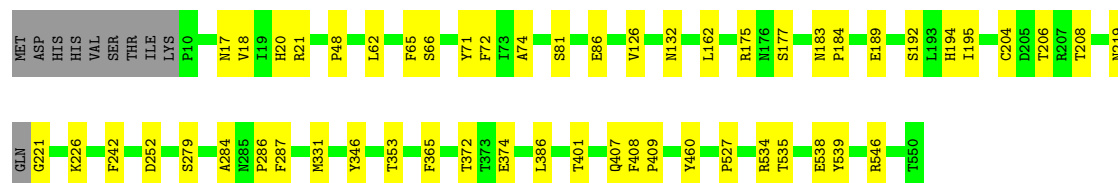
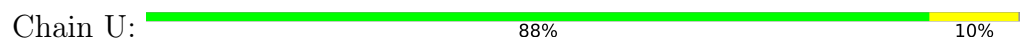
Chain T: 68% 5% 27%



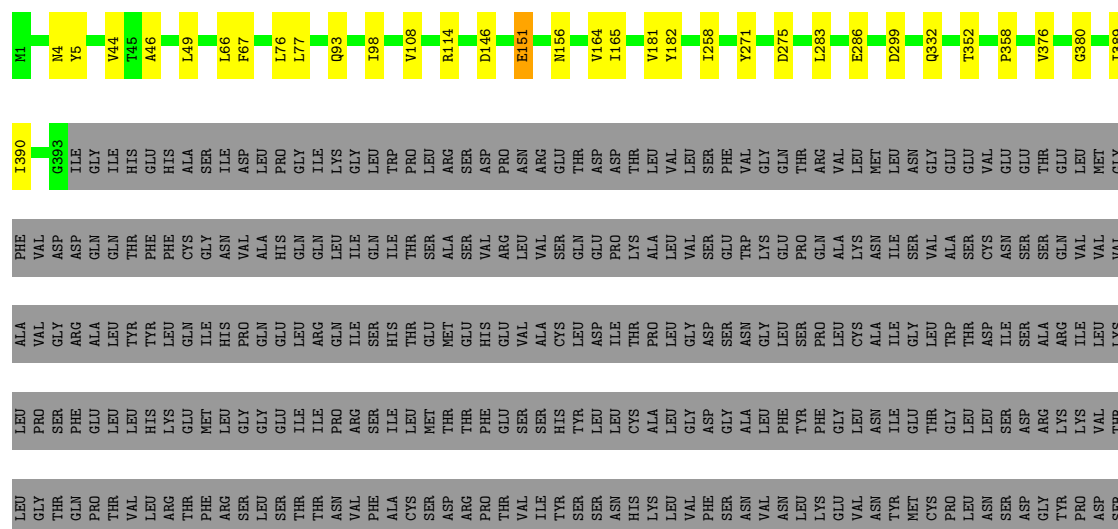
- Molecule 3: DET1 homolog

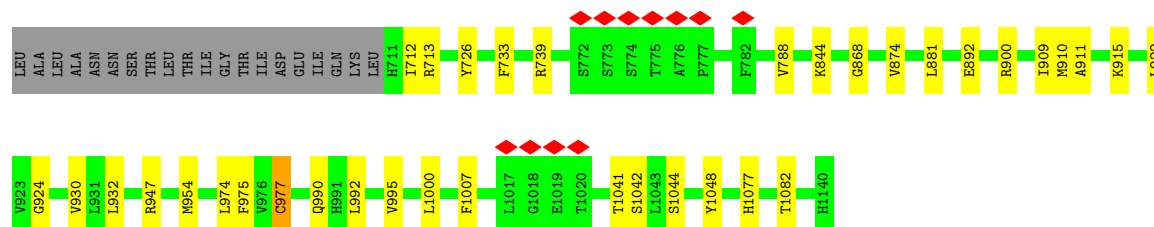


- Molecule 3: DET1 homolog

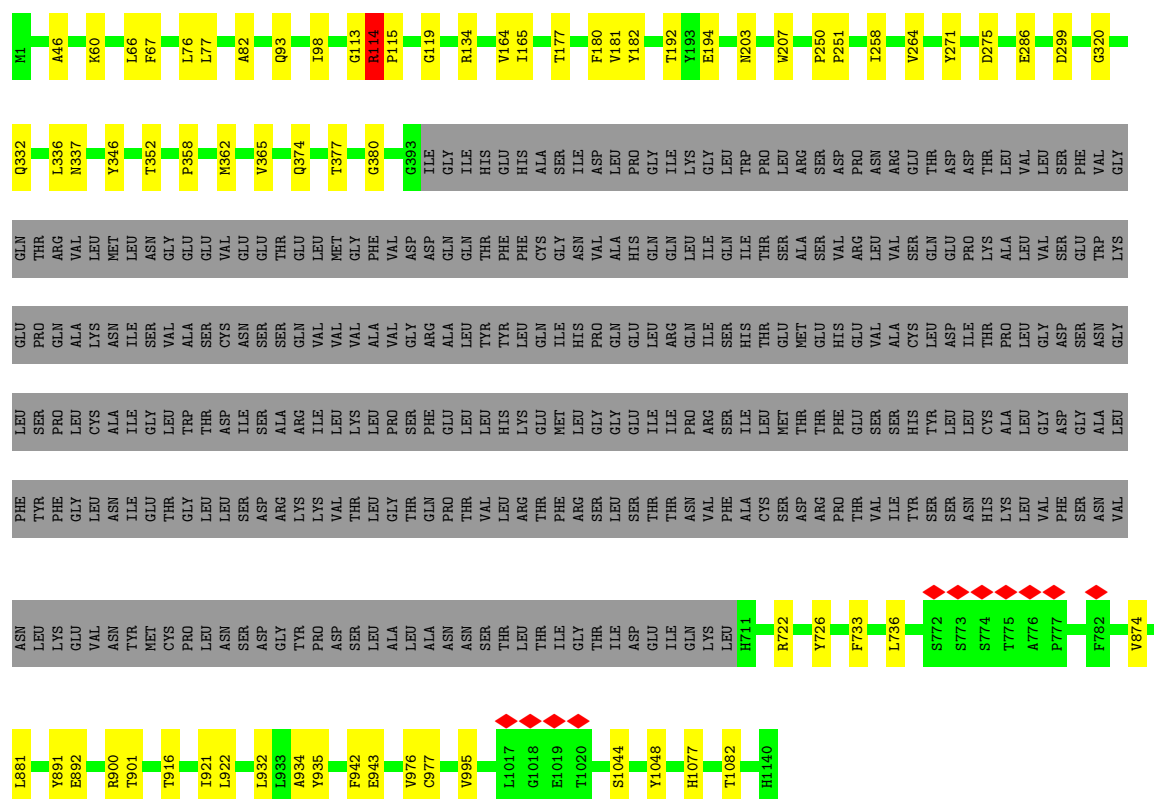


- Molecule 4: DNA damage-binding protein 1

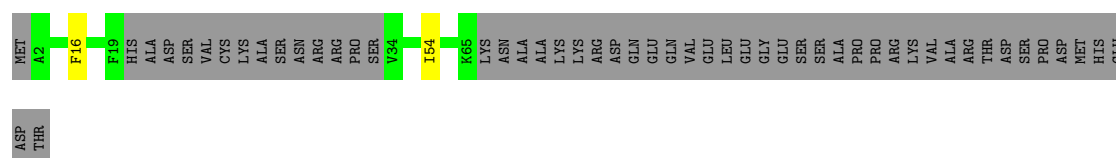




• Molecule 4: DNA damage-binding protein 1



• Molecule 5: DET1- and DDB1-associated protein 1



• Molecule 5: DET1- and DDB1-associated protein 1



HIS	MET
GLU	ALA
ASP	ASP
THR	F4
	F16
	F19
	HIS
	ALA
	ASP
	SER
	VAL
	CYS
	LYS
	ALA
	SER
	ASN
	ARG
	ARG
	PRO
	SER
	V34
	Y35
	L36
	P37
	R66
	ASN
	ALA
	LYS
	LYS
	ARG
	ASP
	GLN
	GLU
	GLN
	VAL
	GLU
	LEU
	GLU
	GLY
	GLU
	SER
	SER
	ALA
	PRO
	PRO
	ARG
	LYS
	VAL
	ALA
	ALA
	ARG
	THR
	ASP
	SER
	PRO
	ASP
	MET

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	72163	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	46.55	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.890	Depositor
Minimum map value	-0.239	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.024	Depositor
Recommended contour level	0.15	Depositor
Map size ($\text{\AA}$)	680.0, 680.0, 680.0	wwPDB
Map dimensions	800, 800, 800	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.85, 0.85, 0.85	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	0.09	0/2224	0.24	0/3100
1	B	0.09	0/2190	0.24	0/3055
1	C	0.09	0/2730	0.26	0/3809
1	D	0.09	0/2525	0.29	0/3525
1	E	0.10	0/2718	0.29	0/3787
1	F	0.09	0/2550	0.23	0/3560
1	G	0.09	0/2178	0.23	0/3039
1	J	0.09	0/2178	0.22	0/3040
2	Q	0.12	0/804	0.28	0/1130
2	T	0.13	0/831	0.31	0/1164
3	R	0.14	0/3036	0.40	0/4214
3	U	0.13	0/3091	0.40	0/4295
4	S	0.13	0/4357	0.33	1/6056 (0.0%)
4	V	0.19	0/4269	0.34	3/5943 (0.1%)
5	j	0.12	0/303	0.29	0/419
5	n	0.09	0/340	0.26	0/469
All	All	0.12	0/36324	0.30	4/50605 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	V	114	ARG	CA-C-N	5.93	125.83	119.85
4	V	114	ARG	C-N-CA	5.93	125.83	119.85
4	V	113	GLY	N-CA-C	5.39	118.45	110.42
4	S	151	GLU	CB-CA-C	-5.18	110.62	116.63

There are no chirality outliers.

There are no planarity outliers.

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	2213	0	1071	16	0
1	B	2182	0	1030	15	0
1	C	2715	0	1286	17	0
1	D	2519	0	1164	19	0
1	E	2704	0	1304	14	0
1	F	2544	0	1174	20	0
1	G	2173	0	1016	25	0
1	J	2173	0	1015	28	0
2	Q	786	0	428	4	0
2	T	811	0	464	5	0
3	R	2976	0	1724	37	0
3	U	3038	0	1769	34	0
4	S	4313	0	2300	39	0
4	V	4234	0	2170	47	0
5	j	299	0	194	2	0
5	n	333	0	258	2	0
All	All	36013	0	18367	312	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 (312) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:R:488:ARG:NH1	3:R:489:PHE:O	2.15	0.79
3:U:409:PRO:HD2	4:V:114:ARG:HB2	1.65	0.79
1:C:425:SER:H	1:C:440:GLY:HA2	1.49	0.78
3:R:21:ARG:HD2	3:R:45:ASN:HD21	1.49	0.76
4:V:119:GLY:O	4:V:134:ARG:NH2	2.21	0.74
1:F:147:TYR:HA	1:F:185:PRO:HA	1.70	0.73
1:F:425:SER:H	1:F:440:GLY:HA2	1.55	0.71
4:S:164:VAL:HA	4:S:182:TYR:HA	1.71	0.71
1:G:659:SER:O	1:G:700:VAL:N	2.24	0.71
1:J:659:SER:O	1:J:700:VAL:N	2.25	0.69
1:E:425:SER:H	1:E:440:GLY:HA2	1.57	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:R:488:ARG:HH12	3:R:497:LEU:HA	1.56	0.69
1:G:435:TYR:HA	1:G:450:TYR:H	1.58	0.69
3:U:408:PHE:CA	4:V:114:ARG:HE	2.10	0.64
1:C:435:TYR:HA	1:C:449:GLU:HA	1.79	0.64
1:F:702:ALA:HB3	1:F:719:ALA:HB3	1.81	0.63
4:V:114:ARG:HG2	4:V:114:ARG:NH1	2.13	0.63
1:D:410:THR:HA	1:D:726:LYS:HA	1.81	0.63
4:S:954:MET:HE1	4:S:975:PHE:HE1	1.64	0.62
1:E:435:TYR:HA	1:E:449:GLU:HA	1.81	0.62
3:U:408:PHE:HA	4:V:114:ARG:HD3	1.82	0.62
4:V:935:TYR:HA	4:V:942:PHE:HA	1.82	0.62
1:E:569:TYR:O	1:E:586:ARG:NH1	2.32	0.62
4:V:114:ARG:HH11	4:V:114:ARG:CG	2.13	0.62
3:U:408:PHE:HA	4:V:114:ARG:HE	1.65	0.62
4:V:332:GLN:HA	4:V:352:THR:HA	1.82	0.62
1:C:498:TRP:HD1	1:C:500:GLY:H	1.46	0.61
1:A:521:PHE:HA	1:A:529:LEU:HA	1.82	0.61
1:J:473:ILE:HA	1:J:490:ASP:HA	1.82	0.61
3:U:409:PRO:CD	4:V:114:ARG:HB2	2.31	0.61
4:V:46:ALA:HA	5:n:16:PHE:H	1.65	0.61
1:D:521:PHE:HA	1:D:529:LEU:HA	1.82	0.61
1:F:410:THR:HA	1:F:726:LYS:HA	1.82	0.61
4:S:165:ILE:N	4:S:181:VAL:O	2.33	0.61
3:U:535:THR:N	3:U:538:GLU:O	2.31	0.61
1:A:407:PRO:HA	1:A:728:LEU:HA	1.83	0.60
3:R:523:HIS:CE1	3:R:528:PHE:HD2	2.19	0.60
3:U:183:ASN:ND2	3:U:353:THR:O	2.33	0.60
1:A:540:LEU:O	1:A:549:VAL:N	2.34	0.60
4:S:46:ALA:HA	5:j:16:PHE:H	1.66	0.60
3:U:372:THR:HG23	3:U:374:GLU:H	1.67	0.60
1:D:700:VAL:HA	1:D:720:ASN:HA	1.85	0.59
4:S:1077:HIS:HA	4:S:1082:THR:HA	1.84	0.59
1:G:425:SER:H	1:G:440:GLY:HA2	1.67	0.59
3:R:345:LYS:HA	3:R:365:PHE:HB3	1.84	0.59
1:F:700:VAL:HA	1:F:720:ASN:HA	1.84	0.59
3:U:401:THR:O	3:U:407:GLN:NE2	2.35	0.59
1:B:521:PHE:HA	1:B:529:LEU:HA	1.84	0.59
1:E:136:CYS:N	1:E:141:ASP:O	2.35	0.59
1:G:624:LEU:O	1:G:633:LEU:N	2.35	0.59
1:J:624:LEU:O	1:J:633:LEU:N	2.36	0.59
4:S:977:CYS:HA	4:S:995:VAL:H	1.68	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:616:ALA:HA	1:G:622:LEU:HA	1.84	0.58
3:R:215:VAL:O	3:R:221:GLY:N	2.36	0.58
3:R:535:THR:N	3:R:538:GLU:O	2.32	0.58
1:D:702:ALA:HB3	1:D:719:ALA:HB3	1.83	0.58
3:U:62:LEU:HA	3:U:74:ALA:HA	1.85	0.58
3:U:242:PHE:HA	3:U:252:ASP:HA	1.86	0.58
1:C:498:TRP:HB3	1:C:505:ARG:HA	1.84	0.58
3:R:379:PHE:HD2	3:R:386:LEU:HD12	1.69	0.58
3:U:18:VAL:HA	3:U:21:ARG:HE	1.68	0.58
1:J:616:ALA:HA	1:J:622:LEU:HA	1.86	0.58
1:B:407:PRO:HA	1:B:728:LEU:HA	1.86	0.58
1:F:521:PHE:HA	1:F:529:LEU:HA	1.85	0.58
4:S:874:VAL:O	4:S:881:LEU:N	2.36	0.57
1:B:540:LEU:O	1:B:549:VAL:N	2.37	0.57
3:R:242:PHE:HA	3:R:252:ASP:HA	1.86	0.57
3:U:65:PHE:HA	3:U:72:PHE:HA	1.86	0.57
1:E:516:CYS:HA	1:E:533:SER:HA	1.87	0.57
1:G:558:VAL:HA	1:G:575:CYS:HA	1.86	0.57
3:R:62:LEU:HA	3:R:74:ALA:HA	1.87	0.57
1:J:583:TYR:HA	1:J:591:PRO:HA	1.86	0.57
3:U:408:PHE:HA	4:V:114:ARG:CD	2.35	0.57
1:C:136:CYS:N	1:C:141:ASP:O	2.39	0.56
1:D:428:GLU:O	1:D:437:ALA:N	2.38	0.56
1:F:428:GLU:O	1:F:437:ALA:N	2.37	0.56
1:C:581:HIS:HA	1:C:594:VAL:HA	1.88	0.56
1:A:666:LEU:O	1:A:675:LEU:N	2.39	0.56
3:U:408:PHE:HA	4:V:114:ARG:NE	2.20	0.56
4:V:358:PRO:HD2	4:V:380:GLY:H	1.70	0.56
1:G:640:ILE:O	1:G:660:GLU:N	2.38	0.56
1:E:640:ILE:O	1:E:660:GLU:N	2.39	0.56
1:D:429:PHE:HA	1:D:436:PHE:HA	1.87	0.55
3:R:520:PHE:HA	3:R:531:SER:HA	1.87	0.55
1:J:666:LEU:O	1:J:675:LEU:N	2.38	0.55
1:E:583:TYR:HA	1:E:592:ILE:H	1.71	0.55
4:S:332:GLN:HA	4:S:352:THR:HA	1.88	0.55
1:G:583:TYR:HA	1:G:591:PRO:HA	1.87	0.55
4:V:66:LEU:HA	4:V:77:LEU:HA	1.88	0.55
1:F:429:PHE:HA	1:F:436:PHE:HA	1.88	0.55
4:S:93:GLN:HA	4:S:98:ILE:HA	1.87	0.55
4:S:358:PRO:HD2	4:S:380:GLY:H	1.72	0.55
1:C:516:CYS:HA	1:C:533:SER:HA	1.88	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:666:LEU:O	1:G:675:LEU:N	2.40	0.55
4:V:114:ARG:HG2	4:V:114:ARG:HH11	1.70	0.55
1:B:666:LEU:O	1:B:675:LEU:N	2.41	0.54
3:R:147:PHE:O	3:R:453:GLN:NE2	2.35	0.54
1:J:640:ILE:O	1:J:660:GLU:N	2.39	0.54
1:J:540:LEU:O	1:J:549:VAL:N	2.40	0.54
1:C:640:ILE:O	1:C:660:GLU:N	2.40	0.54
1:G:407:PRO:HA	1:G:728:LEU:HA	1.90	0.54
1:C:436:PHE:N	1:C:448:TYR:O	2.40	0.54
3:U:331:MET:HA	3:U:346:TYR:HA	1.90	0.54
1:C:429:PHE:HA	1:C:436:PHE:HA	1.90	0.54
3:U:162:LEU:H	3:U:189:GLU:HA	1.73	0.54
4:S:66:LEU:HA	4:S:77:LEU:HA	1.89	0.53
1:G:700:VAL:HA	1:G:720:ASN:HA	1.90	0.53
1:C:135:VAL:HA	1:C:142:MET:HA	1.89	0.53
1:D:447:VAL:O	1:D:464:GLU:N	2.42	0.53
1:C:541:TRP:HD1	1:C:548:SER:HA	1.73	0.53
2:Q:91:TYR:HA	2:Q:188:THR:HG21	1.91	0.53
1:C:437:ALA:HA	1:C:447:VAL:HA	1.90	0.53
1:G:473:ILE:HA	1:G:490:ASP:HA	1.91	0.53
1:D:427:ILE:HA	1:D:438:ILE:HA	1.89	0.53
1:J:476:ILE:HA	1:J:488:SER:HA	1.89	0.53
4:V:114:ARG:NH1	4:V:114:ARG:CG	2.70	0.52
4:V:337:ASN:N	4:V:346:TYR:O	2.42	0.52
1:D:625:TRP:HA	1:D:633:LEU:H	1.73	0.52
1:J:435:TYR:HA	1:J:450:TYR:H	1.73	0.52
3:R:162:LEU:H	3:R:189:GLU:HA	1.75	0.52
4:S:892:GLU:O	4:S:900:ARG:N	2.40	0.52
3:U:86:GLU:N	3:U:126:VAL:O	2.43	0.52
3:U:177:SER:HA	3:U:287:PHE:HA	1.92	0.52
3:R:65:PHE:HA	3:R:72:PHE:HA	1.91	0.52
4:V:93:GLN:HA	4:V:98:ILE:HA	1.92	0.52
1:A:425:SER:H	1:A:440:GLY:HA2	1.75	0.51
1:G:540:LEU:O	1:G:549:VAL:N	2.43	0.51
3:U:279:SER:HA	3:U:284:ALA:HB2	1.91	0.51
1:B:541:TRP:HA	1:B:548:SER:HA	1.91	0.51
1:G:520:ASP:N	1:G:530:ALA:O	2.43	0.51
1:D:411:LEU:N	1:D:725:ILE:O	2.31	0.51
1:J:601:ALA:O	1:J:618:THR:N	2.44	0.51
2:T:122:CYS:O	2:T:165:LEU:N	2.43	0.51
1:J:425:SER:H	1:J:440:GLY:HA2	1.76	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:U:534:ARG:HA	3:U:539:TYR:HA	1.92	0.51
1:J:579:CYS:HA	1:J:596:LYS:HA	1.93	0.51
4:S:910:MET:HG3	4:S:911:ALA:N	2.26	0.51
1:D:478:TRP:HA	1:D:486:LEU:HA	1.93	0.50
3:U:386:LEU:O	3:U:386:LEU:HD23	2.12	0.50
1:A:530:ALA:HA	1:A:540:LEU:HA	1.92	0.50
1:F:427:ILE:HA	1:F:438:ILE:HA	1.93	0.50
1:F:447:VAL:O	1:F:464:GLU:N	2.44	0.50
3:R:534:ARG:HA	3:R:539:TYR:HA	1.92	0.50
1:A:624:LEU:O	1:A:633:LEU:N	2.44	0.50
3:R:333:LYS:O	3:R:345:LYS:N	2.38	0.50
1:E:538:VAL:O	1:E:552:ILE:N	2.42	0.50
1:J:474:SER:N	1:J:489:SER:O	2.44	0.50
4:V:726:TYR:HA	4:V:733:PHE:HA	1.93	0.50
4:V:874:VAL:O	4:V:881:LEU:N	2.34	0.50
1:D:702:ALA:O	1:D:719:ALA:N	2.40	0.50
1:E:581:HIS:HA	1:E:594:VAL:HA	1.92	0.50
3:R:86:GLU:N	3:R:126:VAL:O	2.42	0.50
1:B:667:TYR:HA	1:B:675:LEU:H	1.77	0.49
1:G:411:LEU:O	1:G:725:ILE:N	2.38	0.49
1:E:135:VAL:HA	1:E:142:MET:HA	1.94	0.49
4:V:165:ILE:N	4:V:181:VAL:O	2.42	0.49
1:A:625:TRP:HA	1:A:633:LEU:H	1.78	0.49
1:D:476:ILE:HA	1:D:488:SER:HA	1.93	0.49
1:J:521:PHE:HA	1:J:529:LEU:HA	1.94	0.49
1:F:667:TYR:HA	1:F:675:LEU:H	1.78	0.49
1:A:656:ALA:HA	1:A:666:LEU:HA	1.95	0.49
3:R:486:PRO:HB3	3:R:502:GLN:HA	1.95	0.49
2:Q:122:CYS:O	2:Q:165:LEU:N	2.46	0.48
3:R:223:TYR:CD2	3:R:336:LEU:HB2	2.47	0.48
3:R:195:ILE:O	3:R:204:CYS:N	2.47	0.48
4:S:258:ILE:HA	4:S:275:ASP:HA	1.95	0.48
1:B:530:ALA:HA	1:B:540:LEU:HA	1.95	0.48
1:E:437:ALA:HA	1:E:447:VAL:HA	1.95	0.48
1:F:625:TRP:HA	1:F:633:LEU:H	1.79	0.48
4:V:192:THR:OG1	4:V:207:TRP:O	2.31	0.48
1:F:170:CYS:N	1:F:175:TYR:O	2.47	0.48
3:U:219:ASN:O	3:U:221:GLY:N	2.46	0.48
1:J:574:GLY:HA2	1:J:580:VAL:HA	1.96	0.48
3:R:194:HIS:HA	3:R:206:THR:HA	1.96	0.48
3:U:175:ARG:HA	3:U:286:PRO:HG2	1.96	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:640:ILE:O	1:D:660:GLU:N	2.36	0.48
4:V:164:VAL:HA	4:V:182:TYR:HA	1.96	0.48
4:V:264:VAL:H	4:V:271:TYR:HA	1.78	0.48
1:A:541:TRP:HA	1:A:548:SER:HA	1.94	0.47
3:R:523:HIS:CE1	3:R:528:PHE:CD2	3.01	0.47
1:F:583:TYR:HA	1:F:591:PRO:HA	1.95	0.47
3:U:192:SER:HA	3:U:208:THR:HA	1.96	0.47
1:F:715:VAL:HA	1:F:729:GLU:HA	1.96	0.47
1:G:663:SER:HA	1:G:679:LYS:HA	1.97	0.47
4:S:947:ARG:O	4:S:992:LEU:N	2.44	0.47
1:G:474:SER:N	1:G:489:SER:O	2.47	0.47
4:V:164:VAL:HG11	4:V:180:PHE:HB2	1.96	0.47
1:A:529:LEU:O	1:A:541:TRP:N	2.40	0.47
4:S:389:ILE:O	4:S:713:ARG:N	2.45	0.47
4:S:947:ARG:N	4:S:990:GLN:O	2.42	0.47
3:U:195:ILE:O	3:U:204:CYS:N	2.47	0.47
1:G:667:TYR:HA	1:G:675:LEU:H	1.80	0.47
4:S:286:GLU:N	4:S:299:ASP:O	2.35	0.47
4:S:726:TYR:HA	4:S:733:PHE:HA	1.96	0.47
4:S:5:TYR:N	4:S:1041:THR:O	2.47	0.47
1:E:486:LEU:O	1:E:498:TRP:N	2.48	0.47
4:V:934:ALA:O	4:V:943:GLU:N	2.40	0.47
4:V:892:GLU:O	4:V:900:ARG:N	2.47	0.46
3:R:85:TYR:HA	3:R:127:LEU:HA	1.96	0.46
4:S:910:MET:HE3	4:S:911:ALA:H	1.81	0.46
4:V:194:GLU:O	4:V:203:ASN:N	2.48	0.46
1:A:665:TYR:HA	1:A:677:THR:HA	1.97	0.46
1:E:408:LEU:N	1:E:727:VAL:O	2.43	0.46
1:F:475:CYS:O	1:F:489:SER:N	2.47	0.46
1:G:601:ALA:O	1:G:618:THR:N	2.48	0.46
1:B:493:GLY:HA2	1:B:516:CYS:H	1.81	0.46
1:J:407:PRO:HA	1:J:728:LEU:HA	1.97	0.46
2:T:68:SER:O	2:T:82:THR:N	2.41	0.46
1:G:476:ILE:HA	1:G:488:SER:HA	1.97	0.46
4:V:286:GLU:N	4:V:299:ASP:O	2.35	0.46
1:C:583:TYR:HA	1:C:592:ILE:H	1.81	0.45
1:J:659:SER:H	1:J:700:VAL:N	2.14	0.45
1:B:580:VAL:O	1:B:595:PHE:N	2.34	0.45
4:S:1044:SER:O	4:S:1048:TYR:N	2.47	0.45
1:A:472:LYS:O	1:A:491:TYR:N	2.49	0.45
1:A:667:TYR:HA	1:A:675:LEU:H	1.82	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:V:916:THR:HA	4:V:921:ILE:HA	1.99	0.45
4:V:976:VAL:O	4:V:995:VAL:N	2.50	0.45
1:C:663:SER:HA	1:C:679:LYS:HA	1.98	0.45
1:F:540:LEU:O	1:F:549:VAL:N	2.50	0.45
1:F:582:TYR:O	1:F:592:ILE:N	2.51	0.45
4:V:977:CYS:HA	4:V:995:VAL:H	1.82	0.45
1:B:497:LEU:O	1:B:506:SER:N	2.39	0.44
1:B:656:ALA:HA	1:B:666:LEU:HA	2.00	0.44
1:G:613:ILE:O	1:G:625:TRP:N	2.50	0.44
4:V:365:VAL:O	4:V:374:GLN:N	2.50	0.44
4:V:1044:SER:O	4:V:1048:TYR:N	2.45	0.44
4:S:4:ASN:HA	4:S:1042:SER:HA	1.99	0.44
4:V:891:TYR:HA	4:V:901:THR:HA	1.99	0.44
1:J:411:LEU:O	1:J:725:ILE:N	2.46	0.44
4:V:722:ARG:N	4:V:736:LEU:O	2.51	0.44
3:R:340:ASN:O	3:R:341:HIS:ND1	2.51	0.44
4:S:67:PHE:N	4:S:76:LEU:O	2.34	0.44
3:R:478:ARG:CB	3:R:479:PRO:HD3	2.48	0.44
1:G:582:TYR:O	1:G:592:ILE:N	2.46	0.44
4:V:258:ILE:HA	4:V:275:ASP:HA	1.99	0.44
3:U:17:ASN:O	3:U:20:HIS:N	2.41	0.43
5:n:35:TYR:CE2	5:n:37:PRO:HG3	2.53	0.43
4:V:114:ARG:HA	4:V:115:PRO:HD3	1.84	0.43
1:J:565:PRO:HD2	1:J:610:GLY:HA2	2.01	0.43
3:R:81:SER:HA	3:R:132:ASN:HA	1.99	0.43
3:R:159:ALA:HA	3:R:191:TYR:HA	2.01	0.43
3:U:194:HIS:HA	3:U:206:THR:HA	2.00	0.43
3:R:17:ASN:O	3:R:21:ARG:NH2	2.51	0.43
1:B:529:LEU:O	1:B:541:TRP:N	2.50	0.43
1:J:663:SER:HA	1:J:679:LYS:HA	1.99	0.43
4:S:844:LYS:HA	4:S:868:GLY:HA2	2.01	0.43
1:J:520:ASP:O	1:J:530:ALA:N	2.52	0.43
3:U:460:TYR:OH	3:U:527:PRO:HB3	2.19	0.43
4:S:271:TYR:N	4:S:283:LEU:O	2.44	0.42
1:C:403:ASN:N	1:C:731:VAL:O	2.42	0.42
3:R:408:PHE:HB3	4:S:114:ARG:NH1	2.34	0.42
3:U:48:PRO:HA	3:U:546:ARG:HA	2.00	0.42
3:R:192:SER:HA	3:R:208:THR:HA	2.01	0.42
4:S:156:ASN:HB2	5:j:54:ILE:HD13	2.00	0.42
4:S:915:LYS:NZ	4:S:1007:PHE:O	2.53	0.42
1:G:413:TYR:N	1:G:723:GLY:O	2.45	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:V:177:THR:HA	4:V:194:GLU:HA	2.01	0.42
1:C:170:CYS:N	1:C:175:TYR:O	2.51	0.42
1:J:582:TYR:O	1:J:592:ILE:N	2.48	0.42
4:S:44:VAL:HA	4:S:49:LEU:HA	2.01	0.42
2:T:181:ASP:C	2:T:185:ARG:HE	2.23	0.42
3:U:66:SER:N	3:U:71:TYR:O	2.51	0.42
4:V:250:PRO:HA	4:V:251:PRO:HD3	1.97	0.42
1:A:497:LEU:O	1:A:506:SER:N	2.39	0.42
4:S:376:VAL:HA	4:S:389:ILE:HA	2.02	0.42
3:U:183:ASN:OD1	3:U:184:PRO:HD2	2.20	0.42
1:B:624:LEU:O	1:B:633:LEU:N	2.53	0.42
1:D:136:CYS:N	1:D:141:ASP:O	2.51	0.42
3:R:459:PRO:HD3	3:R:524:PRO:O	2.19	0.42
4:V:60:LYS:HB3	4:V:82:ALA:HB2	2.02	0.42
3:R:74:ALA:O	3:R:83:GLU:N	2.44	0.42
3:R:408:PHE:HB3	4:S:114:ARG:CZ	2.50	0.42
4:S:390:ILE:HA	4:S:712:ILE:HA	2.02	0.42
1:D:583:TYR:HA	1:D:591:PRO:HA	2.02	0.41
1:D:624:LEU:O	1:D:633:LEU:N	2.53	0.41
1:J:495:VAL:N	1:J:509:TYR:O	2.41	0.41
3:R:365:PHE:HE2	3:R:386:LEU:HD13	1.85	0.41
4:S:739:ARG:N	4:S:788:VAL:O	2.42	0.41
1:G:397:SER:O	1:G:401:ARG:N	2.54	0.41
1:J:522:ASN:N	1:J:528:LEU:O	2.46	0.41
1:J:715:VAL:HA	1:J:729:GLU:HA	2.03	0.41
3:R:48:PRO:HA	3:R:546:ARG:HA	2.02	0.41
4:V:362:MET:HA	4:V:377:THR:HA	2.03	0.41
2:Q:68:SER:O	2:Q:82:THR:N	2.39	0.41
4:S:974:LEU:HD11	4:S:1000:LEU:HD11	2.02	0.41
3:U:81:SER:HA	3:U:132:ASN:HA	2.02	0.41
1:E:474:SER:N	1:E:489:SER:O	2.51	0.41
1:G:579:CYS:HA	1:G:596:LYS:HA	2.02	0.41
3:R:143:GLU:O	3:R:218:HIS:N	2.54	0.41
4:V:922:LEU:HA	4:V:932:LEU:HA	2.03	0.41
1:B:665:TYR:HA	1:B:677:THR:HA	2.02	0.41
1:F:136:CYS:N	1:F:141:ASP:O	2.34	0.41
1:F:435:TYR:HA	1:F:449:GLU:HA	2.03	0.41
2:Q:110:PRO:HA	2:Q:111:PRO:HD3	1.94	0.41
4:S:108:VAL:HG12	4:S:108:VAL:O	2.20	0.41
1:A:570:HIS:HA	1:A:584:ASP:HA	2.02	0.41
1:B:493:GLY:HA2	1:B:516:CYS:N	2.36	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:S:146:ASP:H	4:S:151:GLU:H	1.67	0.41
4:S:922:LEU:HA	4:S:932:LEU:HA	2.03	0.41
4:S:924:GLY:HA2	4:S:930:VAL:HA	2.03	0.41
3:U:365:PHE:HE2	3:U:386:LEU:HD13	1.86	0.41
2:T:69:ALA:HA	2:T:81:SER:HA	2.03	0.40
2:T:86:PRO:HD2	2:T:91:TYR:O	2.21	0.40
4:V:320:GLY:O	4:V:336:LEU:N	2.41	0.40
1:D:425:SER:H	1:D:440:GLY:HA2	1.86	0.40
1:J:410:THR:HA	1:J:726:LYS:HA	2.04	0.40
1:D:717:ILE:HA	1:D:727:VAL:HA	2.04	0.40
4:V:67:PHE:O	4:V:76:LEU:N	2.55	0.40
4:V:1077:HIS:HA	4:V:1082:THR:HA	2.02	0.40
1:J:705:TRP:HA	1:J:716:LEU:HA	2.04	0.40
3:R:358:ASP:HA	3:R:359:PRO:HD3	1.97	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	428/731 (58%)	417 (97%)	11 (3%)	0	100	100
1	B	428/731 (58%)	418 (98%)	10 (2%)	0	100	100
1	C	524/731 (72%)	515 (98%)	9 (2%)	0	100	100
1	D	493/731 (67%)	483 (98%)	9 (2%)	1 (0%)	43	77
1	E	519/731 (71%)	505 (97%)	14 (3%)	0	100	100
1	F	498/731 (68%)	491 (99%)	7 (1%)	0	100	100
1	G	429/731 (59%)	423 (99%)	6 (1%)	0	100	100
1	J	429/731 (59%)	417 (97%)	12 (3%)	0	100	100
2	Q	145/201 (72%)	142 (98%)	3 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	T	145/201 (72%)	143 (99%)	2 (1%)	0	100	100
3	R	518/550 (94%)	504 (97%)	12 (2%)	2 (0%)	30	66
3	U	536/550 (98%)	512 (96%)	23 (4%)	1 (0%)	43	77
4	S	819/1140 (72%)	790 (96%)	27 (3%)	2 (0%)	43	77
4	V	819/1140 (72%)	796 (97%)	23 (3%)	0	100	100
5	j	46/102 (45%)	43 (94%)	3 (6%)	0	100	100
5	n	45/102 (44%)	45 (100%)	0	0	100	100
All	All	6821/9834 (69%)	6644 (97%)	171 (2%)	6 (0%)	49	83

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	247	GLN
3	R	226	LYS
4	S	909	ILE
3	R	478	ARG
4	S	977	CYS
3	U	226	LYS

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	15/634 (2%)	15 (100%)	0	100	100
1	B	10/634 (2%)	10 (100%)	0	100	100
1	C	19/634 (3%)	19 (100%)	0	100	100
1	D	11/634 (2%)	11 (100%)	0	100	100
1	E	22/634 (4%)	22 (100%)	0	100	100
1	F	11/634 (2%)	11 (100%)	0	100	100
1	G	8/634 (1%)	8 (100%)	0	100	100
1	J	8/634 (1%)	8 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	Q	18/177 (10%)	18 (100%)	0	100	100
2	T	22/177 (12%)	22 (100%)	0	100	100
3	R	89/499 (18%)	89 (100%)	0	100	100
3	U	91/499 (18%)	91 (100%)	0	100	100
4	S	71/999 (7%)	71 (100%)	0	100	100
4	V	53/999 (5%)	52 (98%)	1 (2%)	50	66
5	j	14/93 (15%)	14 (100%)	0	100	100
5	n	25/93 (27%)	25 (100%)	0	100	100
All	All	487/8608 (6%)	486 (100%)	1 (0%)	85	85

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	V	114	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	512	HIS
3	R	485	HIS
3	R	523	HIS
4	S	163	HIS
3	U	415	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 ⓘ

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.

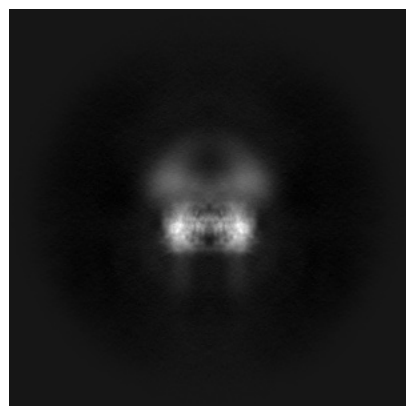
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-63565. These allow visual inspection of the internal detail of the map and identification of artifacts.

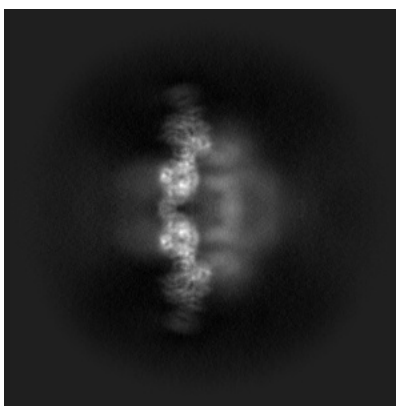
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

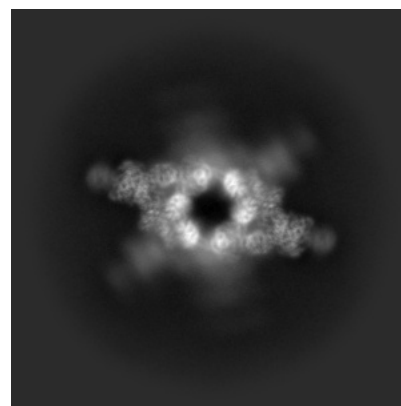
6.1.1 Primary map



X

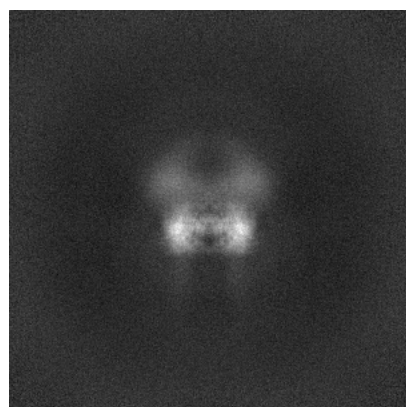


Y

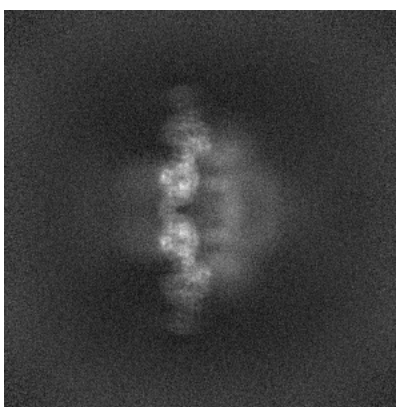


Z

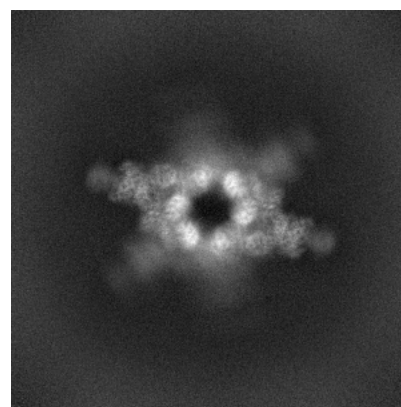
6.1.2 Raw map



X



Y

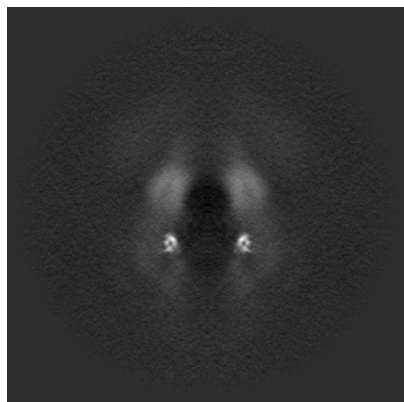


Z

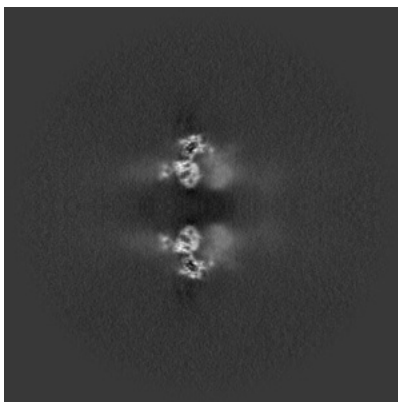
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

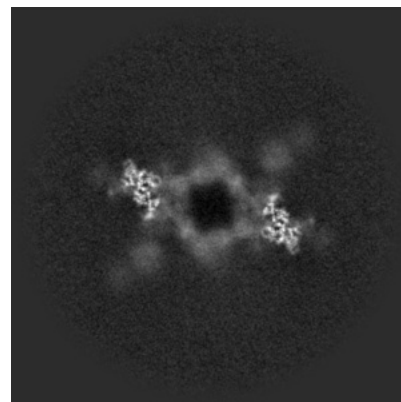
6.2.1 Primary map



X Index: 400

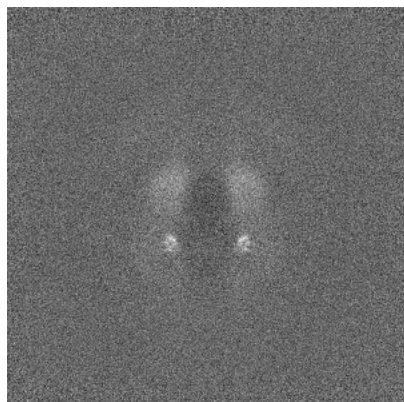


Y Index: 400

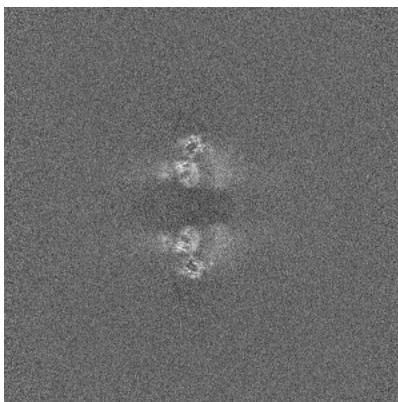


Z Index: 400

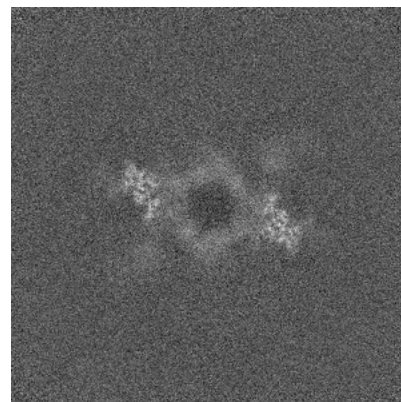
6.2.2 Raw map



X Index: 400



Y Index: 400

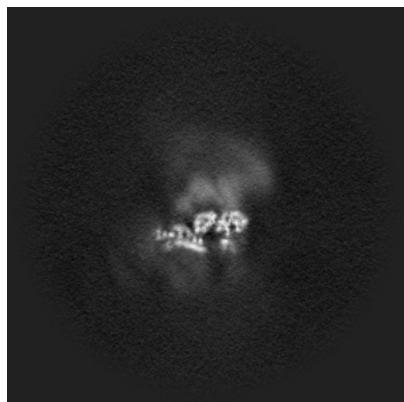


Z Index: 400

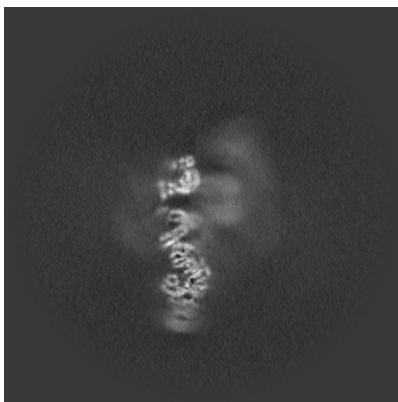
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

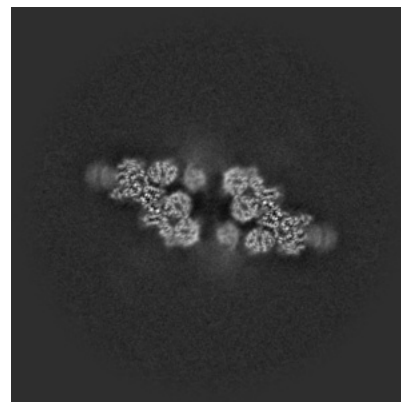
6.3.1 Primary map



X Index: 455

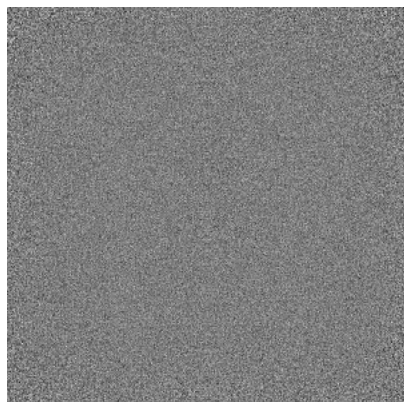


Y Index: 458

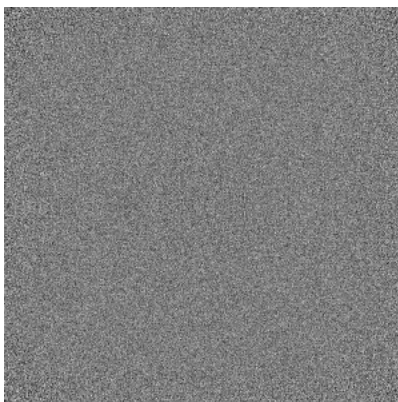


Z Index: 365

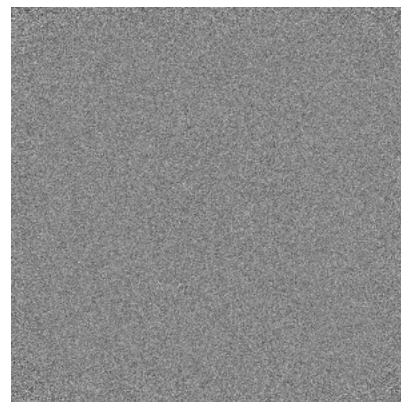
6.3.2 Raw map



X Index: 0



Y Index: 0

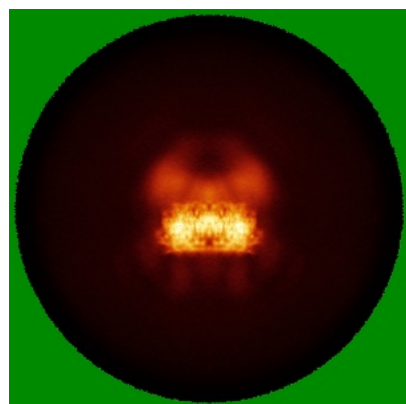


Z Index: 0

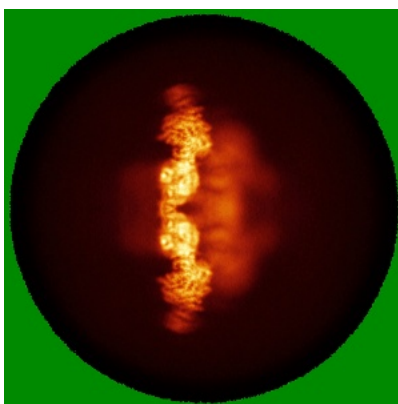
The images above show the largest variance slices of the map in three orthogonal directions.

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

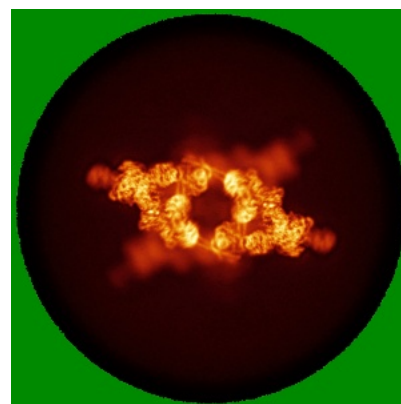
6.4.1 Primary map



X

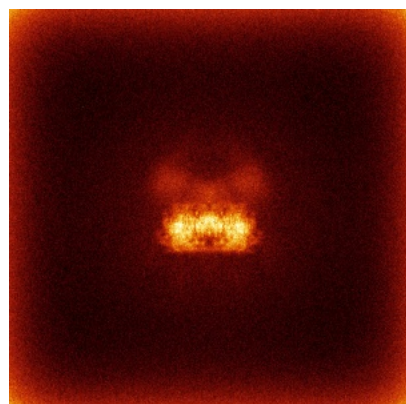


Y

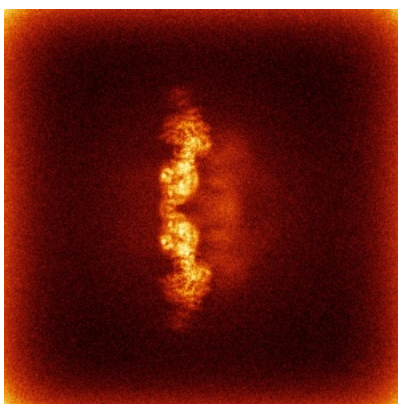


Z

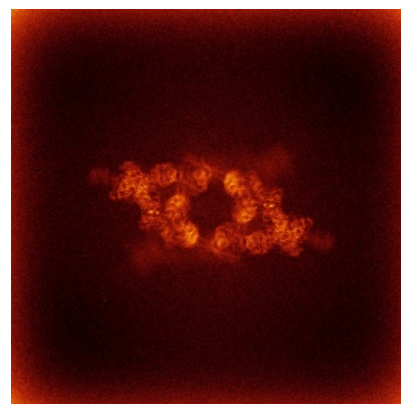
6.4.2 Raw map



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.

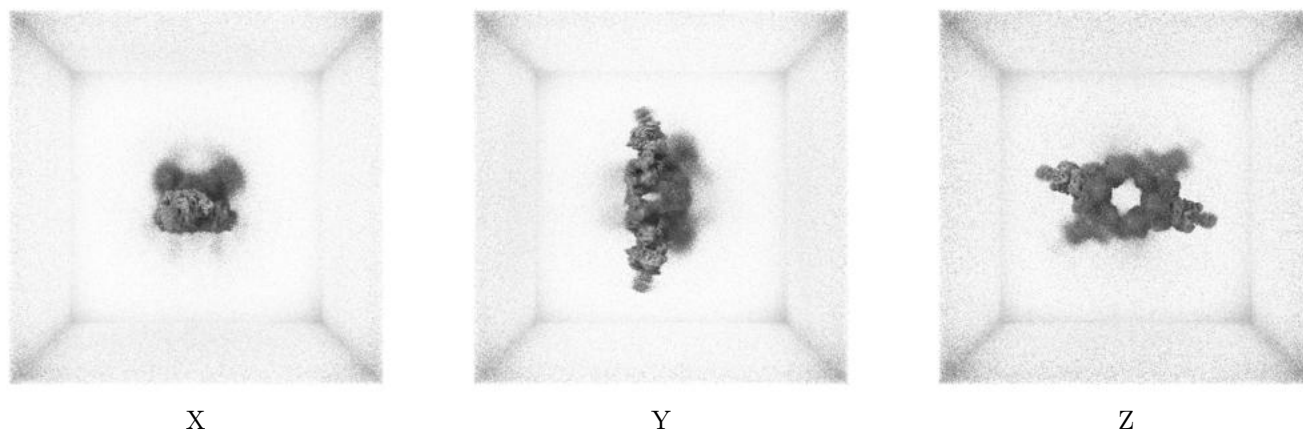
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.15. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

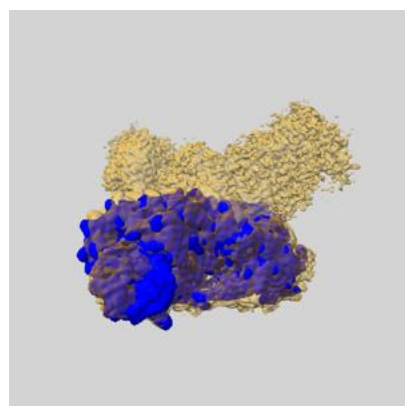
6.6 Mask visualisation [i](#)

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

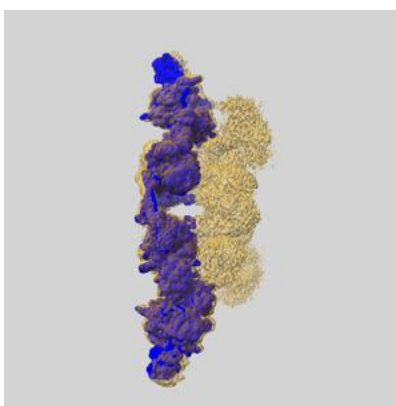
A mask typically either:

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

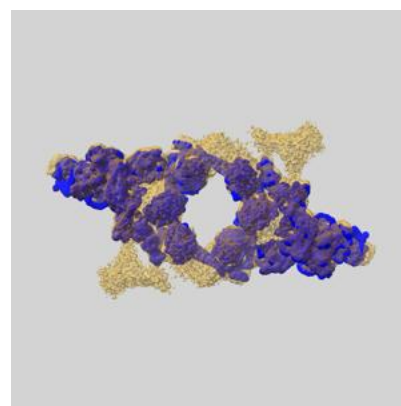
6.6.1 emd_63565_msk_1.map [i](#)



X



Y

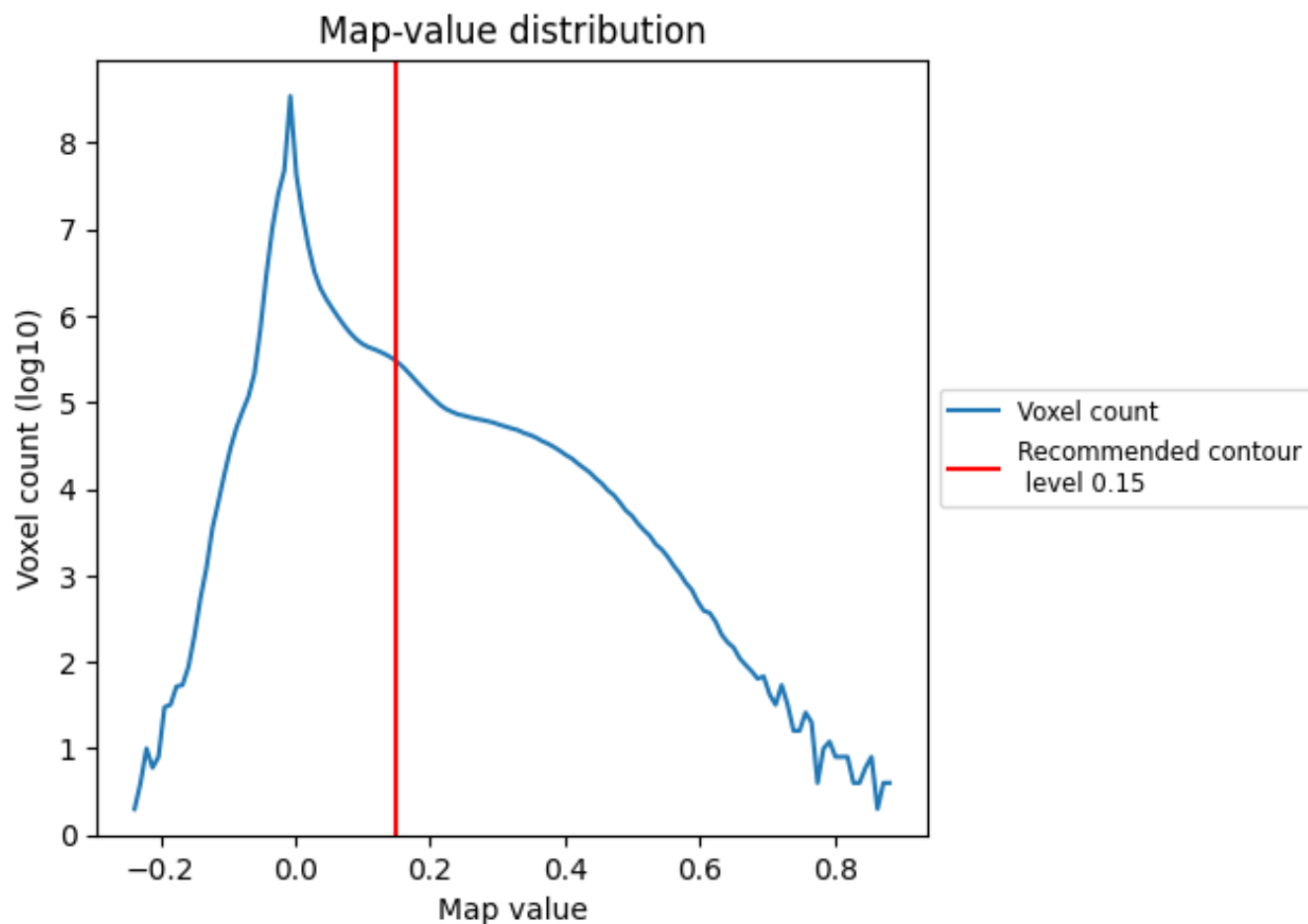


Z

7 Map analysis [i](#)

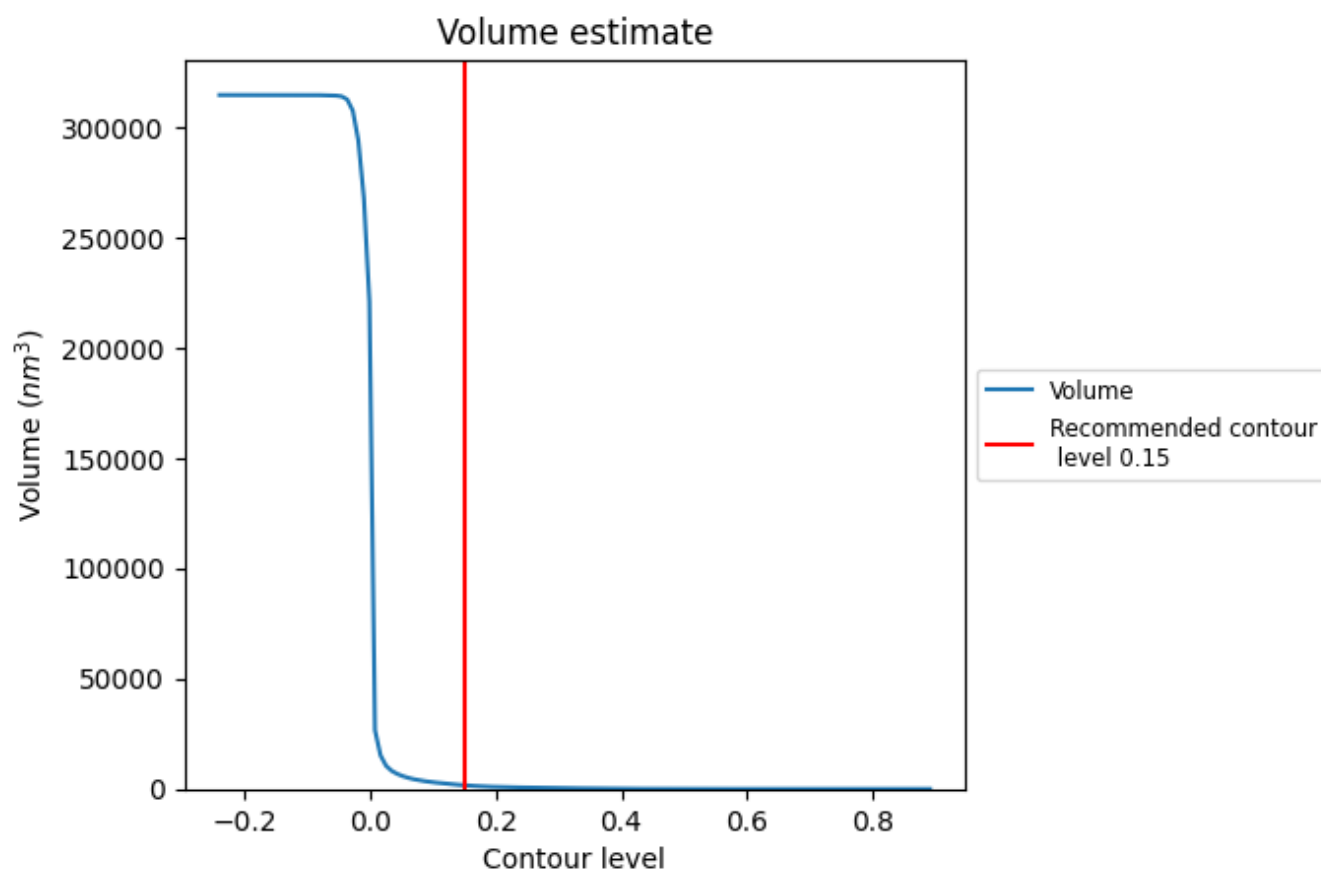
This section contains the results of statistical analysis of the map.

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.

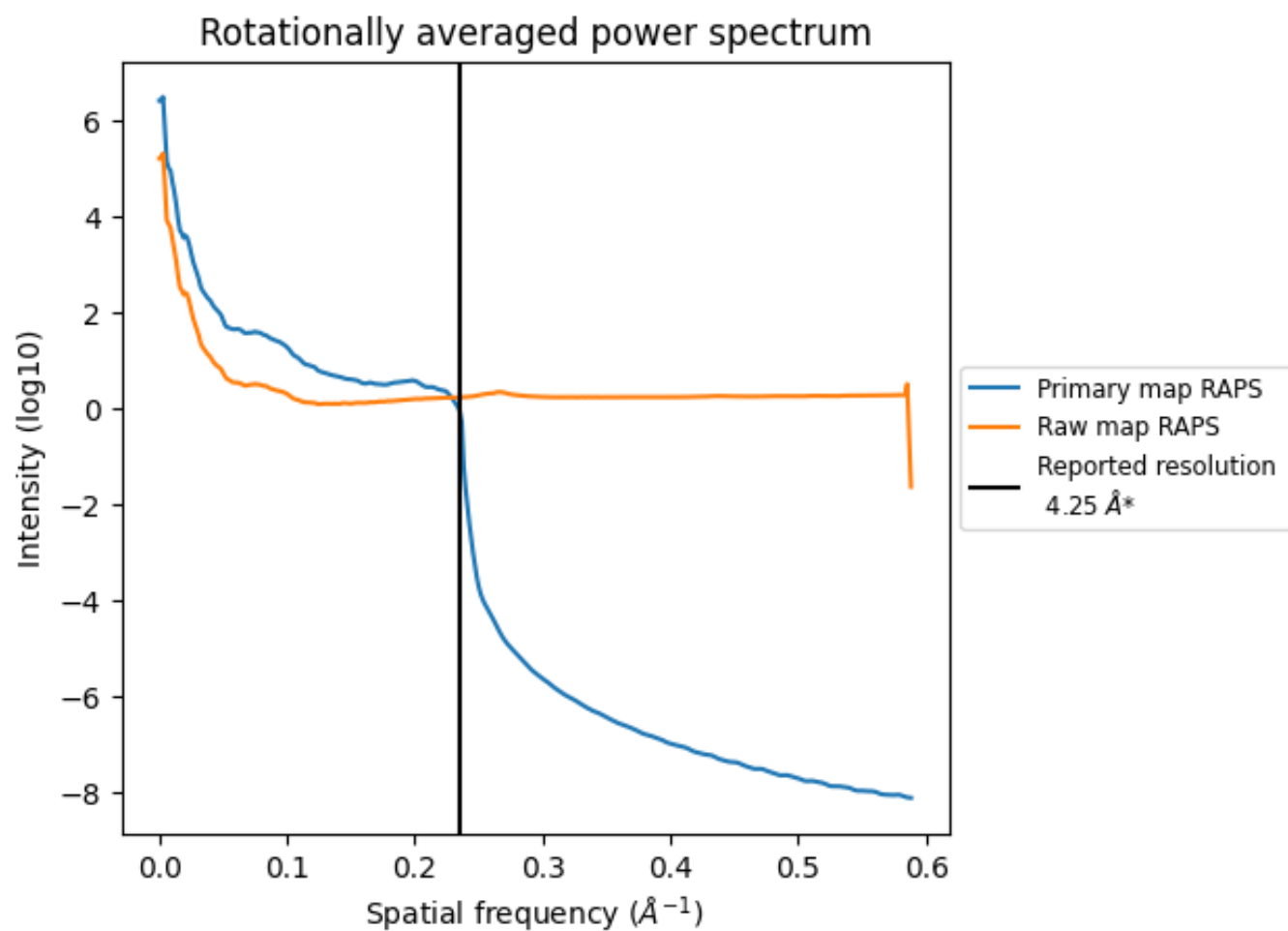
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1680 nm^3 ; this corresponds to an approximate mass of 1518 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 ⓘ

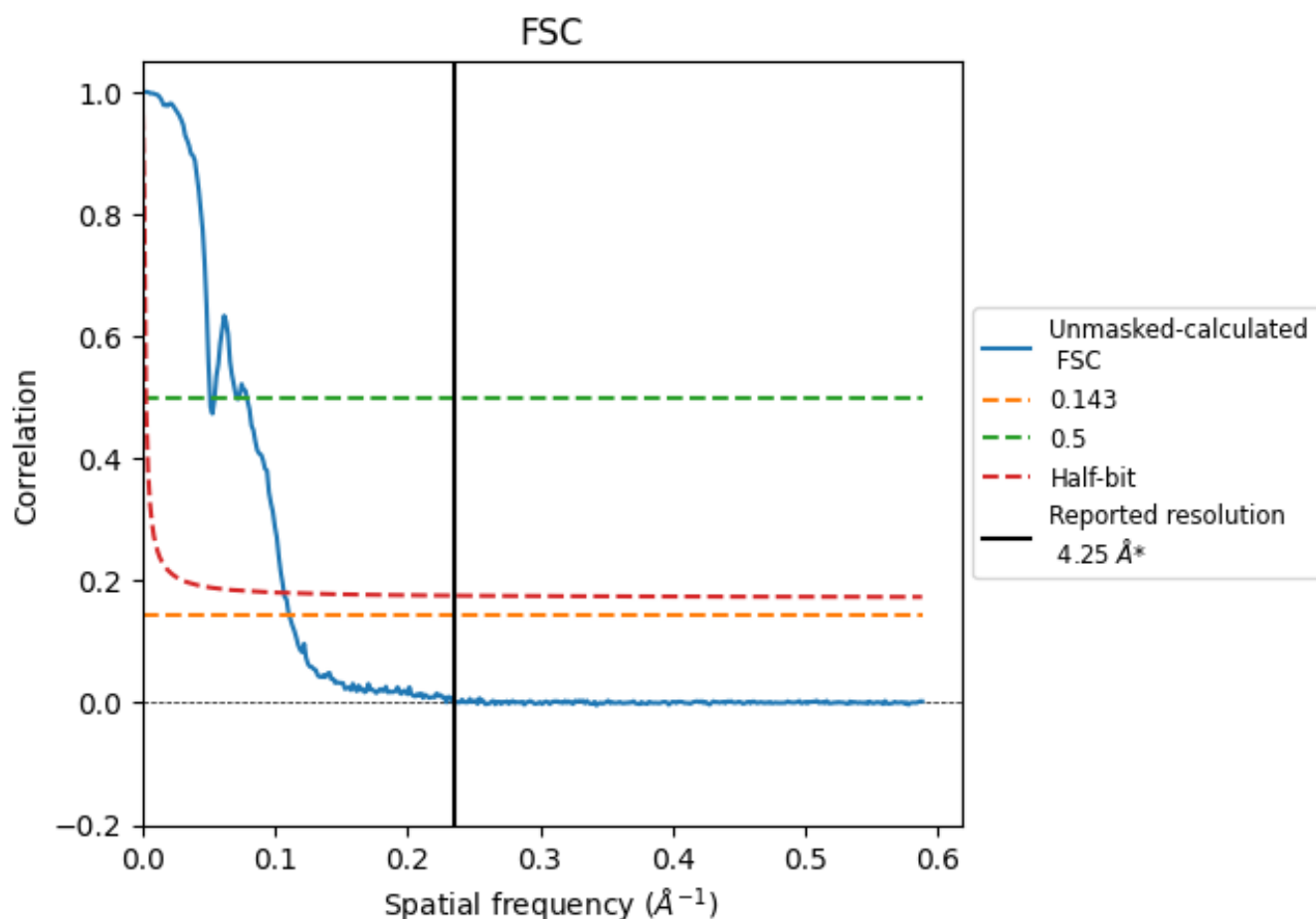


*Reported resolution corresponds to spatial frequency of $0.235 \text{ \AA}^{-1}$

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 0.235 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.25	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	9.02	19.57	9.35

*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 9.02 differs from the reported value 4.25 by more than 10 %

9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-63565 and PDB model 9M0Y. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

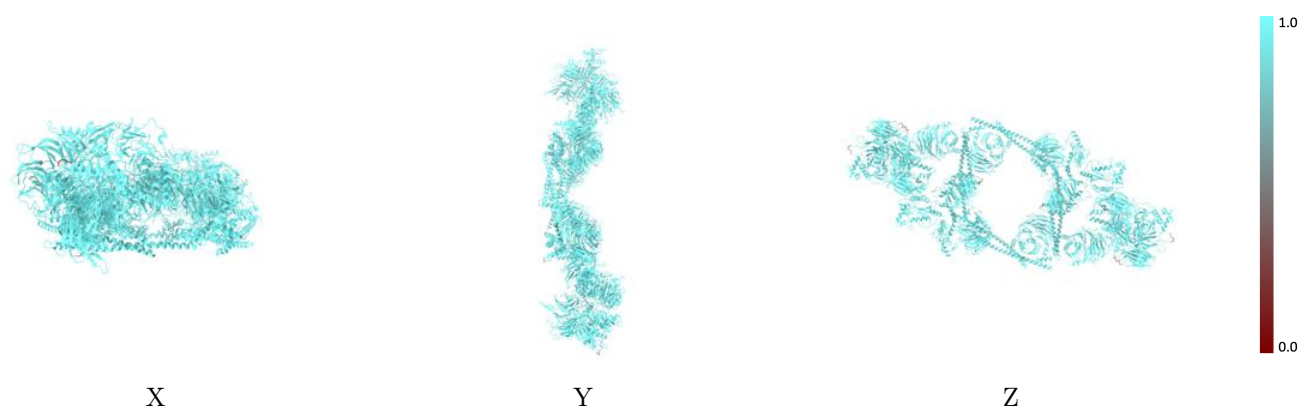
This section was not generated.

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



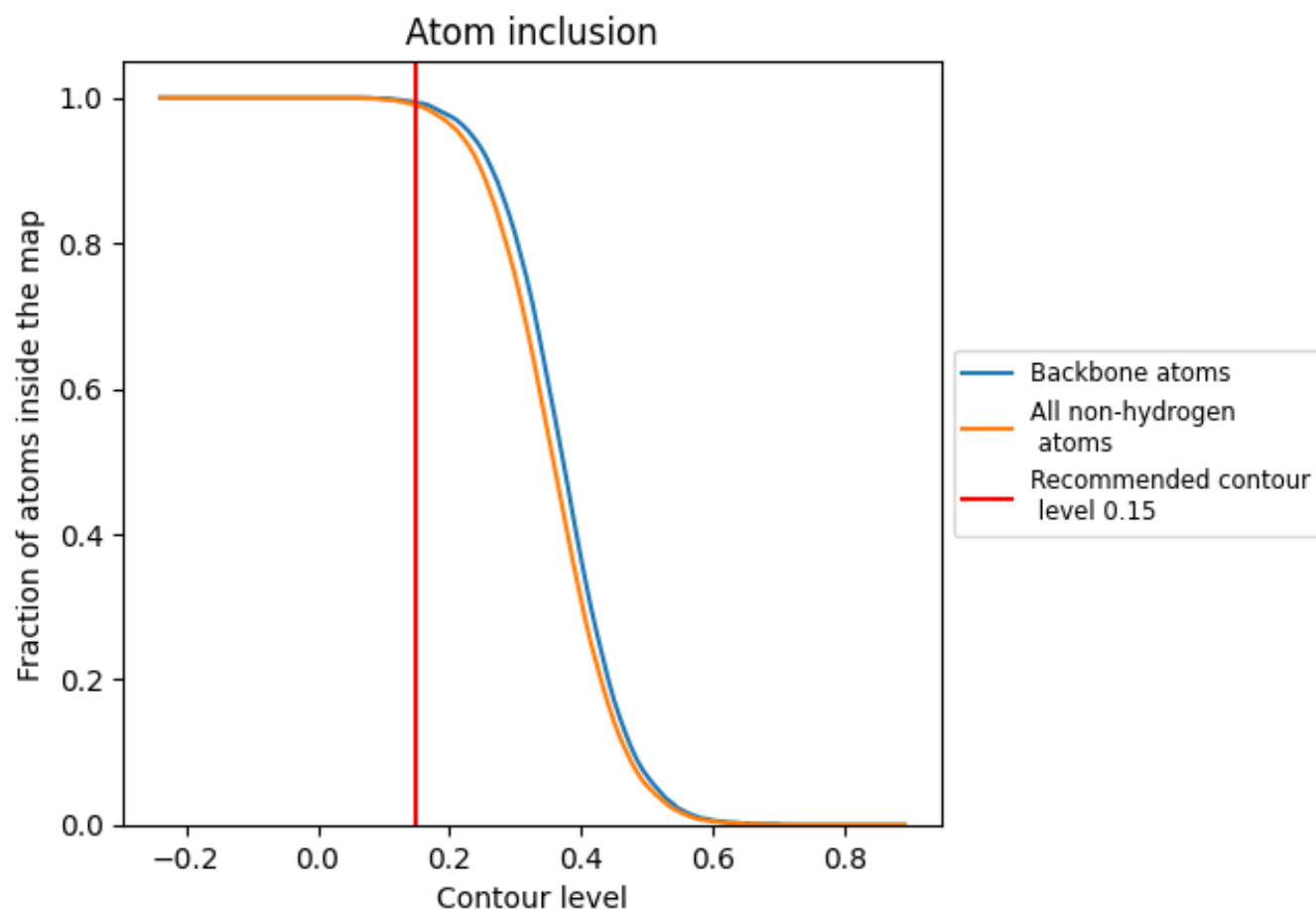
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.15).

9.4 Atom inclusion [i](#)



At the recommended contour level, 99% of all backbone atoms, 99% 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 (0.15) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9900	 0.3220
A	 0.9910	 0.3190
B	 0.9870	 0.3250
C	 0.9930	 0.3240
D	 0.9970	 0.2630
E	 0.9940	 0.3280
F	 0.9970	 0.2540
G	 0.9890	 0.2400
J	 0.9900	 0.2410
Q	 0.9940	 0.3600
R	 0.9940	 0.3690
S	 0.9810	 0.3610
T	 0.9950	 0.3620
U	 0.9920	 0.3640
V	 0.9810	 0.3600
j	 0.9800	 0.3490
n	 0.9880	 0.3220

