



Full wwPDB EM Validation Report ⓘ

Apr 5, 2026 – 05:18 PM UTC

PDB ID : 9PRC / pdb_00009prc
EMDB ID : EMD-71807
Title : HDAG complex with 86-pRNA, Body1
Authors : Itskanov, S.; Lansdon, E.B.
Deposited on : 2025-07-23
Resolution : 3.50 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

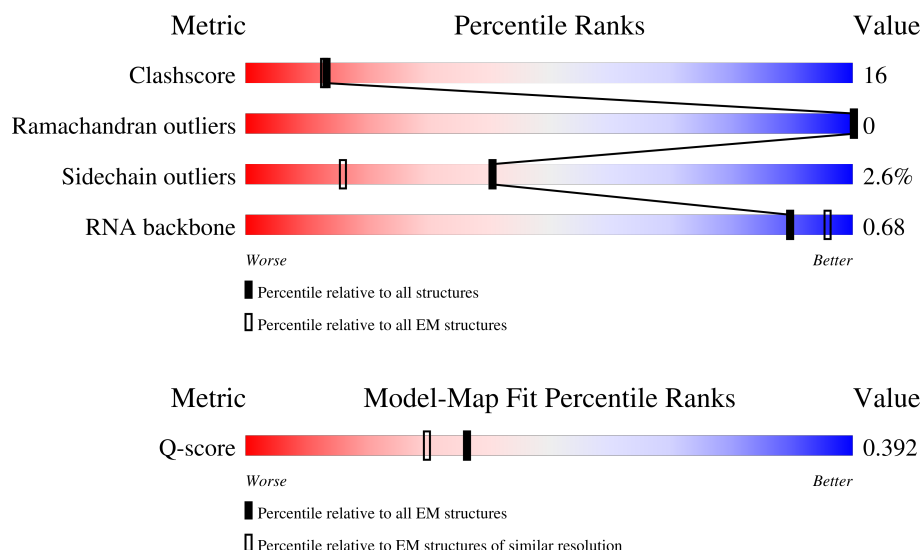
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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.



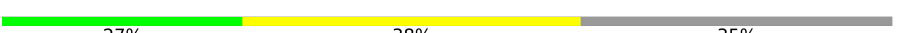
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	13950 (3.00 - 4.00)

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	194	
1	B	194	
1	C	194	

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Mol	Chain	Length	Quality of chain
1	D	194	 16% 9% 75%
1	E	194	 21% • • 74%
1	F	194	 17% 8% 75%
1	G	194	 15% 10% • 74%
1	H	194	 18% 7% • 75%
2	J	55	 27% 38% 35%
2	K	55	 36% 36% 27%
2	M	55	 33% 33% 35%
2	P	55	 5% 44% 24% 27%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6493 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 Small delta antigen.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	52	Total	C	N	O	0	0
			422	268	79	75		
1	B	52	Total	C	N	O	0	0
			422	268	79	75		
1	C	50	Total	C	N	O	0	0
			412	262	77	73		
1	D	49	Total	C	N	O	0	0
			407	259	76	72		
1	E	50	Total	C	N	O	0	0
			412	262	77	73		
1	F	49	Total	C	N	O	0	0
			407	259	76	72		
1	G	50	Total	C	N	O	0	0
			412	262	77	73		
1	H	49	Total	C	N	O	0	0
			407	259	76	72		

- Molecule 2 is a RNA chain called Palindromic RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	J	36	Total	C	N	O	P	0	0
			756	342	126	252	36		
2	K	40	Total	C	N	O	P	0	0
			840	380	140	280	40		
2	M	36	Total	C	N	O	P	0	0
			756	342	126	252	36		
2	P	40	Total	C	N	O	P	0	0
			840	380	140	280	40		

Chain D:  16% 9% 75%

GLU	ARG	ARG	ILE	ALA	GLY	PRO	GLN	VAL	GLY	GLY	VAL	ASN	PRO	LEU	GLY	GLY	THR	ARG	ALA	PRO	GLY	GLY	PHE	VAL	SER	PRO	MET	GLN	VAL	GLU	GLY	SER	PRO	GLY	ILE	ASP	GLY	GLN	PHE	PRO										
VAL	ASP	SER	GLY	PRO	GLY	LYS	ARG	PRO	VAL	SER	ARG	GLY	PHE	THR	ASP	LYS	GLU	ARG	HIS	ARG	ARG	GLY	LYS	ALA	LEU	GLU	ASN	ASN	GLN	LEU	SER	ALA	GLY	PRO	THR	ARG	LYS	THR	THR	GLY	ILE	ASP	GLY	GLN	PHE	PRO				
SER	ARG	PRO	GLY	GLY	ARG	LYS	ASN	ARG	GLY	GLY	R13	E14	L17	K25	K26	E29	L30	L34	T41	L44	E47	H48	P49	W50	L51	K55	G56	I57	L58	G59	K60	LYS	ASP	LYS	ASP	GLY	GLY	GLY	ALA	PRO	PRO	ALA	LYS	ARG	ALA	ALA	THR	ASP	GLN	MET

• Molecule 1: Small delta antigen

Chain E:  21% . . 74%

SER	ARG	PRO	GLU	GLY	ARG	LYS	ASN	ARG	GLY	G12	V16	L17	K26	K36	I41	H48	P49	W50	L51	I57	L58	K61	ASP	LYS	GLY	GLY	GLY	GLY	ALA	PRO	PRO	LYS	ARG	LYS	ALA	ARG	THR	ASP	SER	GLY	ILE	ARG	PRO	LYS	ARG	PRO	GLY	ARG
GLY	THR	ASP	GLY	LYS	ARG	GLN	ARG	ASP	HIS	ARG	ARG	ARG	LYS	ALA	LEU	ASN	LYS	GLN	LEU	SER	ALA	GLY	LEU	SER	LYS	GLY	GLU	GLY	ASP	GLY	ILE	ARG	GLY	THR	GLY	ASP	GLY	ASP	GLY	ARG	GLY	THR	GLY	VAL	GLN	VAL	GLY	GLY
VAL	ASN	LEU	GLY	GLY	THR	ARG	GLY	ALA	PRO	GLY	GLY	PHE	VAL	PRO	SER	MET	GLY	VAL	PRO	GLU	SER	PRO	GLY	GLY	GLY	GLY	GLY	LEU	ILE	ARG	SER	GLN	PHE	PRO	GLY	THR	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY

• Molecule 1: Small delta antigen

Chain F:  17% 8% 75%

GLU	ARG	ARG	ILE	ALA	GLY	PRO	GLN	VAL	GLY	GLY	VAL	ASN	PRO	LEU	GLY	GLY	THR	ARG	ALA	PRO	GLY	PHE	VAL	SER	PRO	MET	GLN	VAL	PRO	GLU	SER	PRO	PHE	THR	ARG	THR	GLY	GLY	GLY	GLY	ILE	GLY	GLY	GLY	GLN	GLY	PHE	PRO																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																													
VAL	ASP	SER	GLY	PRO	ARG	LYS	ARG	PRO	SER	ARG	GLY	GLY	PHE	THR	ASP	LYS	GLU	ARG	GLN	HIS	ARG	ARG	LYS	ALA	LEU	GLU	ASN	LYS	ARG	GLY	ALA	GLY	LYS	GLN	LEU	SER	ALA	GLY	GLY	LYS	ASN	LEU	SER	LYS	LYS	GLY	GLY	ALA	ALA	GLY	GLY	ASP	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY</

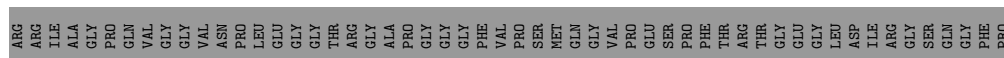
• Molecule 1: Small delta antigen

Chain G:  15% 10% . 74%

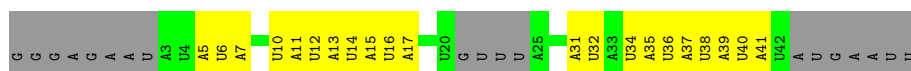
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GLN	MET	GLU	VAL	ASP	SER	GLY	PRO	ARG	GLY	ARG	PRO	SER	ARG	GLY	GLY	PHE	THR	LYS	GLU	ARG	GLN	ASP	HIS	ARG	ARG	LYS	ALA	LEU	ASN	LYS	LEU	SER	ALA	GLY	LYS	ASN	LEU	SER	LYS	GLU	GLU	GLU	GLU	GLU	GLU	GLU	GLU		
GLU	ARG	ARG	GLU	ARG	ARG	ILE	ALA	ALA	GLY	GLN	VAL	GLY	VAL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

• Molecule 1: Small delta antigen

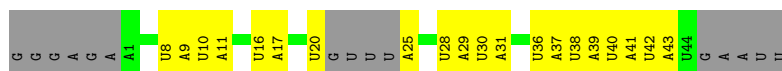
Chain H:  18% 7% . 75%

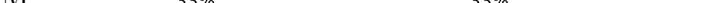


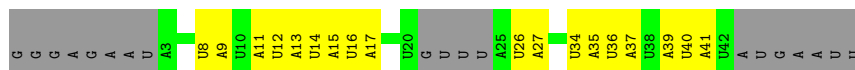
Chain J:  27% 38% 35%

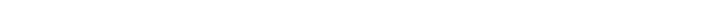


Chain K: 36% 36% 27%



Chain M:  33% 33% 35%



Chain P:  5% 44% 24% 27%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	298422	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	0.012	Depositor
Minimum map value	-0.006	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.000	Depositor
Recommended contour level	0.00115	Depositor
Map size (Å)	349.91998, 349.91998, 349.91998	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.729, 0.729, 0.729	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	0.18	0/427	0.38	0/567
1	B	0.16	0/427	0.30	0/567
1	C	0.19	0/417	0.38	0/553
1	D	0.19	0/412	0.36	0/546
1	E	0.19	0/417	0.38	0/553
1	F	0.21	0/412	0.46	0/546
1	G	0.18	0/417	0.32	0/553
1	H	0.20	0/412	0.37	0/546
2	J	0.14	0/844	0.19	0/1306
2	K	0.11	0/938	0.21	0/1452
2	M	0.14	0/844	0.19	0/1306
2	P	0.15	0/938	0.33	0/1452
All	All	0.16	0/6905	0.31	0/9947

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	422	0	437	17	0
1	B	422	0	437	20	0
1	C	412	0	433	14	0
1	D	407	0	431	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	412	0	433	13	0
1	F	407	0	431	19	0
1	G	412	0	433	25	0
1	H	407	0	431	17	0
2	J	756	0	380	19	0
2	K	840	0	422	22	0
2	M	756	0	380	18	0
2	P	840	0	422	32	0
All	All	6493	0	5070	185	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 16.

All (185) 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:F:26:LYS:HA	1:F:26:LYS:HE2	1.64	0.77
1:E:26:LYS:HA	1:E:26:LYS:HE2	1.69	0.74
2:P:4:U:H3	2:P:41:A:H2	1.38	0.69
1:B:31:GLU:OE2	1:B:35:ARG:NH1	2.25	0.69
1:A:24:ARG:HH21	1:H:14:GLU:HG3	1.59	0.67
2:M:14:U:H2'	2:M:15:A:H8	1.58	0.67
1:G:23:GLY:HA3	1:H:41:ILE:HD11	1.77	0.65
1:B:29:GLU:OE2	1:B:32:ARG:NH1	2.29	0.65
1:D:55:LYS:HA	1:D:58:LEU:HD12	1.78	0.65
2:M:14:U:H2'	2:M:15:A:C8	2.32	0.65
2:K:38:U:H2'	2:K:39:A:H8	1.64	0.63
1:F:12:GLY:N	1:F:15:GLU:OE2	2.33	0.62
1:G:55:LYS:HA	1:G:58:LEU:HD12	1.82	0.62
2:J:36:U:H2'	2:J:37:A:C8	2.37	0.60
2:M:36:U:H2'	2:M:37:A:C8	2.37	0.60
1:G:20:TRP:CD2	1:H:51:LEU:HD11	2.36	0.60
1:B:32:ARG:NH2	2:K:8:U:O3'	2.32	0.60
1:D:13:ARG:HD3	1:F:58:LEU:HA	1.82	0.59
1:G:34:LEU:HG	1:H:30:LEU:HD23	1.84	0.59
1:C:58:LEU:HD21	1:E:17:LEU:HD13	1.82	0.59
2:K:36:U:H2'	2:K:37:A:H8	1.67	0.59
2:K:10:U:H2'	2:K:11:A:H8	1.66	0.59
2:M:12:U:H2'	2:M:13:A:H8	1.68	0.59
2:P:2:U:H2'	2:P:3:A:H8	1.68	0.59
1:F:15:GLU:OE1	1:F:15:GLU:N	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:16:VAL:HG11	1:D:48:HIS:ND1	2.18	0.58
2:P:28:U:H2'	2:P:29:A:C8	2.39	0.58
2:K:42:U:H2'	2:K:43:A:H8	1.69	0.57
2:K:28:U:H2'	2:K:29:A:C8	2.40	0.57
1:C:16:VAL:HG11	1:D:48:HIS:CE1	2.40	0.56
2:J:14:U:H2'	2:J:15:A:C8	2.40	0.56
2:K:28:U:H2'	2:K:29:A:H8	1.69	0.56
2:P:6:U:H2'	2:P:7:A:H8	1.71	0.56
2:K:37:A:H2'	2:K:38:U:H6	1.70	0.56
1:D:17:LEU:HD22	1:F:58:LEU:HD21	1.88	0.56
2:M:8:U:H2'	2:M:9:A:C8	2.41	0.56
1:A:13:ARG:HD3	1:G:58:LEU:HA	1.87	0.55
1:G:48:HIS:HE1	1:H:16:VAL:HG11	1.70	0.55
1:B:54:ILE:HD13	1:G:54:ILE:HG12	1.88	0.55
2:P:28:U:H2'	2:P:29:A:H8	1.72	0.55
2:K:38:U:H2'	2:K:39:A:C8	2.41	0.54
1:F:35:ARG:O	1:F:39:LYS:HG2	2.07	0.54
1:E:58:LEU:HD21	1:G:17:LEU:HD22	1.88	0.54
2:J:5:A:N6	2:J:41:A:N1	2.56	0.54
2:J:6:U:H2'	2:J:7:A:C8	2.42	0.54
1:G:13:ARG:HH21	1:H:50:TRP:CD1	2.25	0.53
1:C:50:TRP:CD1	1:D:13:ARG:HH21	2.27	0.53
2:K:42:U:H2'	2:K:43:A:C8	2.44	0.53
1:G:45:GLU:OE2	1:H:20:TRP:NE1	2.32	0.53
1:B:58:LEU:HD21	1:H:14:GLU:HG2	1.91	0.52
2:M:12:U:H2'	2:M:13:A:C8	2.44	0.52
2:P:16:U:H2'	2:P:17:A:C8	2.44	0.52
1:A:16:VAL:HG11	1:B:48:HIS:ND1	2.25	0.52
1:D:48:HIS:HB3	1:D:50:TRP:NE1	2.25	0.52
2:P:8:U:H2'	2:P:9:A:C8	2.45	0.52
2:P:29:A:H2'	2:P:30:U:C6	2.44	0.52
2:K:37:A:H2'	2:K:38:U:C6	2.45	0.51
2:J:14:U:H2'	2:J:15:A:H8	1.77	0.50
1:F:17:LEU:HD22	1:G:21:VAL:HG21	1.93	0.50
2:P:12:U:HO2'	2:P:13:A:H8	1.57	0.50
1:C:20:TRP:HB2	1:D:44:LEU:HD23	1.94	0.50
2:K:36:U:H2'	2:K:37:A:C8	2.46	0.50
2:P:6:U:H2'	2:P:7:A:C8	2.46	0.50
1:G:12:GLY:N	1:G:15:GLU:OE2	2.45	0.50
1:A:21:VAL:HG21	1:H:17:LEU:HD22	1.93	0.49
2:J:38:U:H2'	2:J:39:A:H8	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:4:U:H2'	2:P:5:A:H8	1.75	0.49
1:F:29:GLU:OE1	1:F:32:ARG:NH2	2.45	0.49
1:E:58:LEU:HD11	1:G:17:LEU:HD23	1.94	0.49
1:A:27:LEU:HD11	1:B:38:LYS:HG3	1.94	0.49
1:G:54:ILE:O	1:G:58:LEU:HG	2.12	0.49
2:J:34:U:H2'	2:J:35:A:H8	1.78	0.49
1:A:17:LEU:HD22	1:G:58:LEU:HD21	1.95	0.49
2:M:39:A:H2'	2:M:40:U:C6	2.48	0.48
2:M:16:U:H2'	2:M:17:A:C8	2.48	0.48
2:M:34:U:H2'	2:M:35:A:H8	1.78	0.48
1:C:40:LYS:HA	1:C:40:LYS:HE2	1.95	0.48
1:D:26:LYS:HA	1:D:26:LYS:HE2	1.95	0.48
1:F:17:LEU:HD13	1:H:58:LEU:HD12	1.96	0.48
1:G:25:LYS:HA	1:G:25:LYS:HD2	1.63	0.48
2:J:38:U:H2'	2:J:39:A:C8	2.48	0.48
1:G:47:GLU:C	1:G:48:HIS:HD2	2.21	0.47
1:A:26:LYS:HG2	1:B:37:VAL:HG21	1.95	0.47
2:P:7:A:H2'	2:P:8:U:C6	2.50	0.47
1:C:48:HIS:HB3	1:C:50:TRP:NE1	2.30	0.47
2:J:31:A:H2'	2:J:32:U:C6	2.49	0.47
2:P:32:U:O2'	2:P:33:A:N3	2.38	0.47
1:C:50:TRP:CH2	1:D:13:ARG:HG3	2.49	0.47
1:G:33:ASP:OD1	1:H:30:LEU:HD21	2.14	0.47
1:E:58:LEU:HA	1:G:13:ARG:HD3	1.96	0.46
1:H:53:ASN:O	1:H:57:ILE:HG22	2.15	0.46
1:A:48:HIS:NE2	1:B:16:VAL:HG21	2.29	0.46
2:P:9:A:H2	2:P:36:U:H3	1.63	0.46
2:M:11:A:H2'	2:M:12:U:C6	2.51	0.46
2:M:16:U:H2'	2:M:17:A:H8	1.81	0.46
2:M:39:A:H2'	2:M:40:U:H6	1.81	0.46
2:J:16:U:H2'	2:J:17:A:C8	2.51	0.46
2:P:11:A:H2'	2:P:12:U:C6	2.51	0.46
2:P:36:U:H2'	2:P:37:A:O4'	2.16	0.46
1:E:36:LYS:HE3	1:E:36:LYS:HA	1.97	0.46
1:G:51:LEU:HD23	1:G:51:LEU:HA	1.79	0.46
2:K:40:U:H2'	2:K:41:A:H8	1.81	0.46
2:K:10:U:H2'	2:K:11:A:C8	2.50	0.45
2:P:2:U:H2'	2:P:3:A:C8	2.49	0.45
1:A:16:VAL:HG21	1:B:48:HIS:CE1	2.51	0.45
1:B:32:ARG:HH21	2:K:9:A:P	2.39	0.45
1:C:23:GLY:HA3	1:D:41:ILE:HD11	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:26:U:H2'	2:M:27:A:H8	1.81	0.45
2:P:3:A:N1	2:P:42:U:O2'	2.39	0.45
1:B:43:LYS:O	1:B:47:GLU:HG2	2.16	0.45
2:M:11:A:H2'	2:M:12:U:H6	1.81	0.45
2:M:40:U:H2'	2:M:41:A:O4'	2.17	0.45
2:J:11:A:H2'	2:J:12:U:H6	1.81	0.45
2:P:40:U:O4	2:P:41:A:N6	2.50	0.45
1:E:50:TRP:CH2	1:F:17:LEU:HD12	2.52	0.44
2:P:19:A:H2'	2:P:19:A:N3	2.33	0.44
1:F:39:LYS:HE2	1:F:39:LYS:HB2	1.84	0.44
2:J:10:U:H2'	2:J:11:A:O4'	2.17	0.44
2:J:31:A:H2'	2:J:32:U:H6	1.82	0.44
1:H:25:LYS:HD3	1:H:25:LYS:HA	1.88	0.44
2:P:38:U:H2'	2:P:39:A:C8	2.52	0.44
1:G:41:ILE:HD11	1:H:23:GLY:C	2.42	0.44
2:P:25:A:H2'	2:P:26:U:O2	2.17	0.44
2:P:27:A:H2'	2:P:28:U:C6	2.53	0.44
1:C:37:VAL:HG21	1:D:26:LYS:HG3	2.00	0.44
1:E:16:VAL:HG11	1:F:48:HIS:CE1	2.53	0.44
1:B:45:GLU:HA	1:B:51:LEU:HD12	2.00	0.44
1:C:30:LEU:HB3	1:D:34:LEU:HG	2.00	0.44
2:K:16:U:H2'	2:K:17:A:C8	2.53	0.44
1:C:33:ASP:HB3	1:D:30:LEU:HD11	2.00	0.44
1:A:45:GLU:OE2	1:B:20:TRP:NE1	2.50	0.43
1:D:25:LYS:HA	1:D:25:LYS:HD2	1.77	0.43
1:D:47:GLU:C	1:D:48:HIS:HD2	2.26	0.43
2:M:34:U:H2'	2:M:35:A:C8	2.53	0.43
1:C:54:ILE:HA	1:C:57:ILE:HG22	1.99	0.43
2:J:12:U:H2'	2:J:13:A:H8	1.84	0.43
2:P:15:A:H2'	2:P:16:U:C6	2.53	0.43
1:G:48:HIS:N	1:G:48:HIS:CD2	2.87	0.43
2:M:8:U:H2'	2:M:9:A:H8	1.84	0.43
1:E:48:HIS:NE2	1:F:16:VAL:HG21	2.33	0.43
2:J:34:U:H2'	2:J:35:A:C8	2.54	0.43
1:A:43:LYS:O	1:A:47:GLU:HG2	2.17	0.43
1:D:34:LEU:HD23	1:D:34:LEU:HA	1.82	0.43
2:P:12:U:O2'	2:P:13:A:H5'	2.18	0.43
1:A:21:VAL:HG11	1:H:18:GLU:HA	2.01	0.43
1:A:50:TRP:CD1	1:B:13:ARG:HH21	2.36	0.43
1:B:39:LYS:NZ	2:K:30:U:H5''	2.33	0.43
2:P:13:A:H2'	2:P:14:U:C6	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:48:HIS:CD2	1:F:16:VAL:HG21	2.54	0.43
2:J:11:A:H2'	2:J:12:U:C6	2.53	0.43
2:P:27:A:H2'	2:P:28:U:H6	1.83	0.43
2:J:35:A:H2'	2:J:36:U:H6	1.84	0.43
1:G:30:LEU:HD23	1:G:30:LEU:HA	1.85	0.42
2:K:40:U:H2'	2:K:41:A:C8	2.54	0.42
2:P:13:A:C4	2:P:32:U:O4	2.72	0.42
2:P:39:A:H2'	2:P:40:U:C6	2.54	0.42
1:E:48:HIS:HB3	1:E:51:LEU:HB2	2.00	0.42
2:J:40:U:H2'	2:J:41:A:C8	2.53	0.42
1:B:26:LYS:HE3	1:B:30:LEU:HD11	2.02	0.42
1:B:44:LEU:HD23	1:B:44:LEU:HA	1.88	0.42
1:G:30:LEU:HD11	1:H:33:ASP:HB3	2.02	0.42
1:B:42:LYS:HB3	1:B:42:LYS:HE3	1.89	0.42
2:J:16:U:H2'	2:J:17:A:H8	1.85	0.41
1:D:48:HIS:CD2	1:D:48:HIS:N	2.86	0.41
1:F:14:GLU:HB2	1:H:58:LEU:HD21	2.03	0.41
2:K:38:U:C2	2:K:39:A:C8	3.08	0.41
2:P:12:U:O2'	2:P:13:A:H8	2.04	0.41
1:D:29:GLU:OE2	1:D:29:GLU:HA	2.20	0.41
1:D:51:LEU:HD23	1:D:51:LEU:HA	1.79	0.41
1:A:20:TRP:NE1	1:B:45:GLU:OE2	2.49	0.41
1:A:44:LEU:HD23	1:A:44:LEU:HA	1.85	0.41
1:C:51:LEU:HD12	1:C:51:LEU:HA	1.71	0.41
1:D:14:GLU:HG2	1:F:58:LEU:HD13	2.01	0.41
1:F:54:ILE:HD13	1:F:54:ILE:HA	1.83	0.41
2:K:25:A:H2'	2:K:25:A:N3	2.34	0.41
2:P:16:U:H2'	2:P:17:A:H8	1.86	0.41
2:P:32:U:H1'	2:P:33:A:O5'	2.20	0.41
1:E:41:ILE:HD11	1:F:23:GLY:C	2.46	0.41
1:E:57:ILE:HD12	1:E:57:ILE:HA	1.93	0.41
2:K:41:A:H2'	2:K:42:U:C6	2.56	0.40
1:F:29:GLU:OE1	1:F:29:GLU:HA	2.21	0.40
2:K:9:A:H2'	2:K:10:U:H6	1.86	0.40
1:A:13:ARG:NH1	1:G:61:LYS:HA	2.36	0.40
1:A:29:GLU:OE2	1:A:29:GLU:HA	2.22	0.40
2:M:26:U:H2'	2:M:27:A:C8	2.57	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	50/194 (26%)	50 (100%)	0	0	100	100
1	B	50/194 (26%)	49 (98%)	1 (2%)	0	100	100
1	C	48/194 (25%)	47 (98%)	1 (2%)	0	100	100
1	D	47/194 (24%)	47 (100%)	0	0	100	100
1	E	48/194 (25%)	47 (98%)	1 (2%)	0	100	100
1	F	47/194 (24%)	47 (100%)	0	0	100	100
1	G	48/194 (25%)	48 (100%)	0	0	100	100
1	H	47/194 (24%)	47 (100%)	0	0	100	100
All	All	385/1552 (25%)	382 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	43/159 (27%)	42 (98%)	1 (2%)	44	64
1	B	43/159 (27%)	43 (100%)	0	100	100
1	C	43/159 (27%)	42 (98%)	1 (2%)	44	64
1	D	43/159 (27%)	42 (98%)	1 (2%)	44	64
1	E	43/159 (27%)	41 (95%)	2 (5%)	23	49
1	F	43/159 (27%)	41 (95%)	2 (5%)	23	49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	43/159 (27%)	42 (98%)	1 (2%)	44	64
1	H	43/159 (27%)	42 (98%)	1 (2%)	44	64
All	All	344/1272 (27%)	335 (97%)	9 (3%)	41	62

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LEU
1	C	38	LYS
1	D	57	ILE
1	E	41	ILE
1	E	51	LEU
1	F	41	ILE
1	F	51	LEU
1	G	15	GLU
1	H	51	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	48	HIS
1	G	48	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	J	34/55 (61%)	0	0
2	K	38/55 (69%)	2 (5%)	0
2	M	34/55 (61%)	0	0
2	P	38/55 (69%)	19 (50%)	1 (2%)
All	All	144/220 (65%)	21 (14%)	1 (0%)

All (21) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	K	20	U
2	K	31	A
2	P	10	U
2	P	11	A

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Mol	Chain	Res	Type
2	P	13	A
2	P	26	U
2	P	27	A
2	P	29	A
2	P	30	U
2	P	31	A
2	P	32	U
2	P	33	A
2	P	34	U
2	P	35	A
2	P	37	A
2	P	38	U
2	P	40	U
2	P	41	A
2	P	42	U
2	P	43	A
2	P	44	U

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	P	32	U

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no 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 ⓘ

There are no chain breaks in this entry.

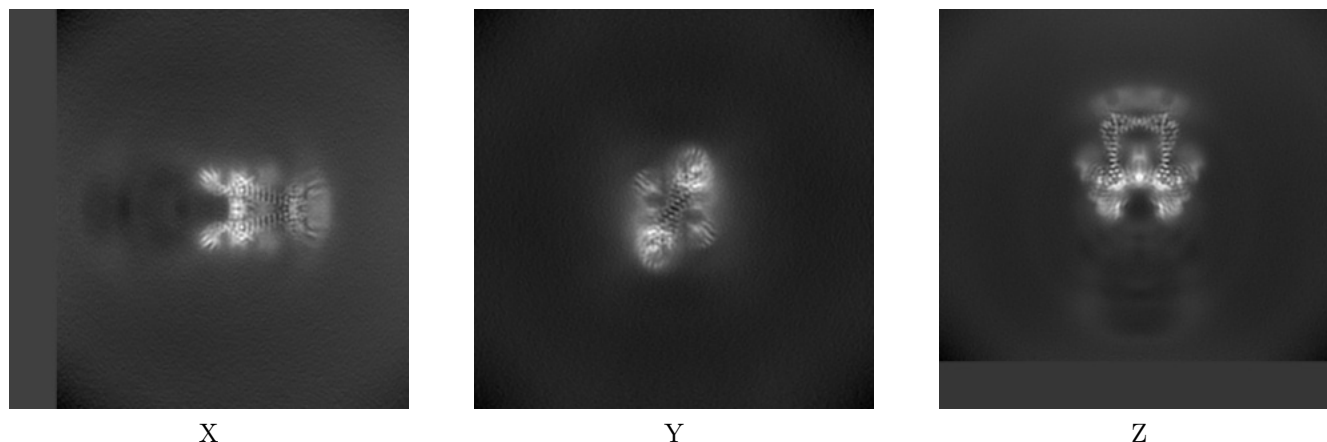
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-71807. These allow visual inspection of the internal detail of the map and identification of artifacts.

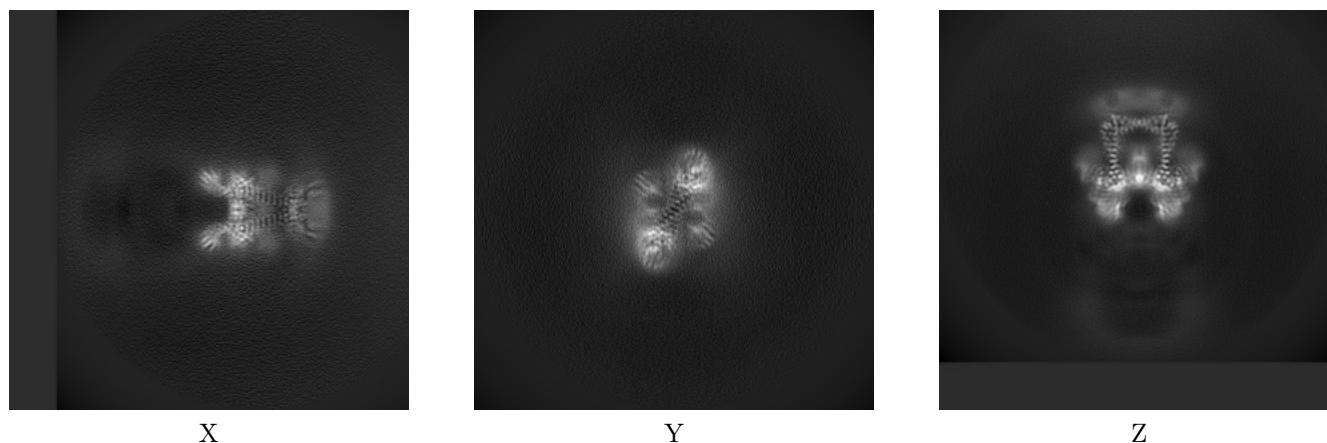
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

6.1.1 Primary map



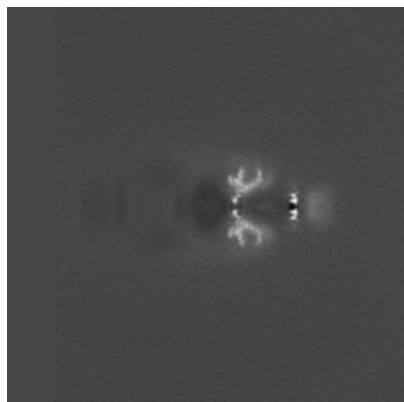
6.1.2 Raw map



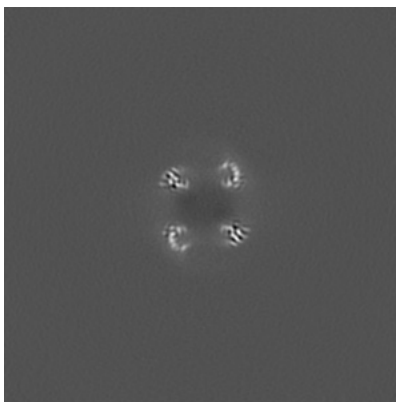
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

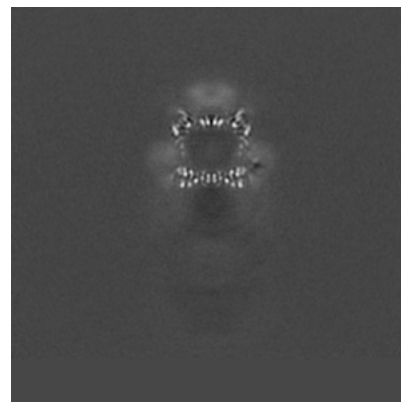
6.2.1 Primary map



X Index: 240

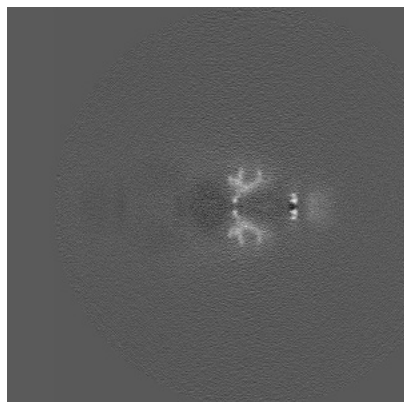


Y Index: 240



Z Index: 240

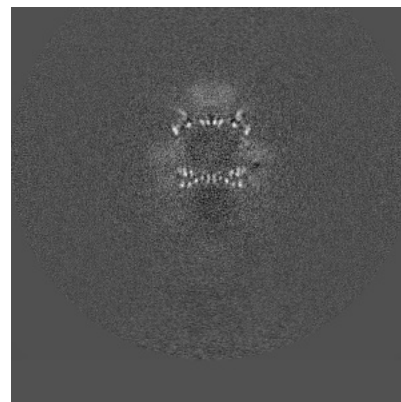
6.2.2 Raw map



X Index: 240



Y Index: 240

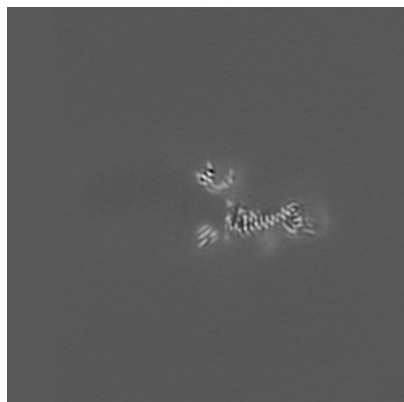


Z Index: 240

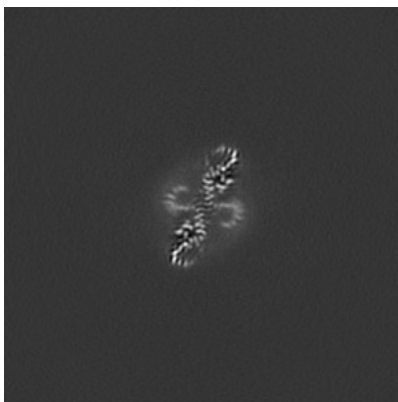
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 212

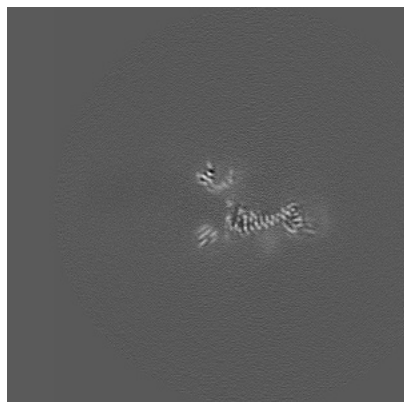


Y Index: 279



Z Index: 263

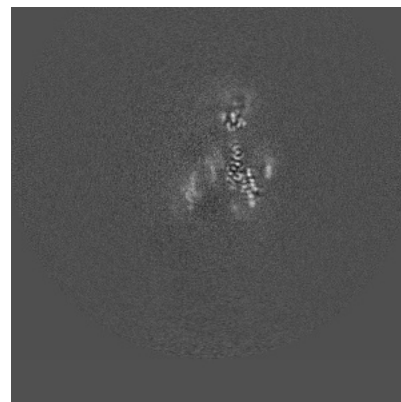
6.3.2 Raw map



X Index: 212



Y Index: 279

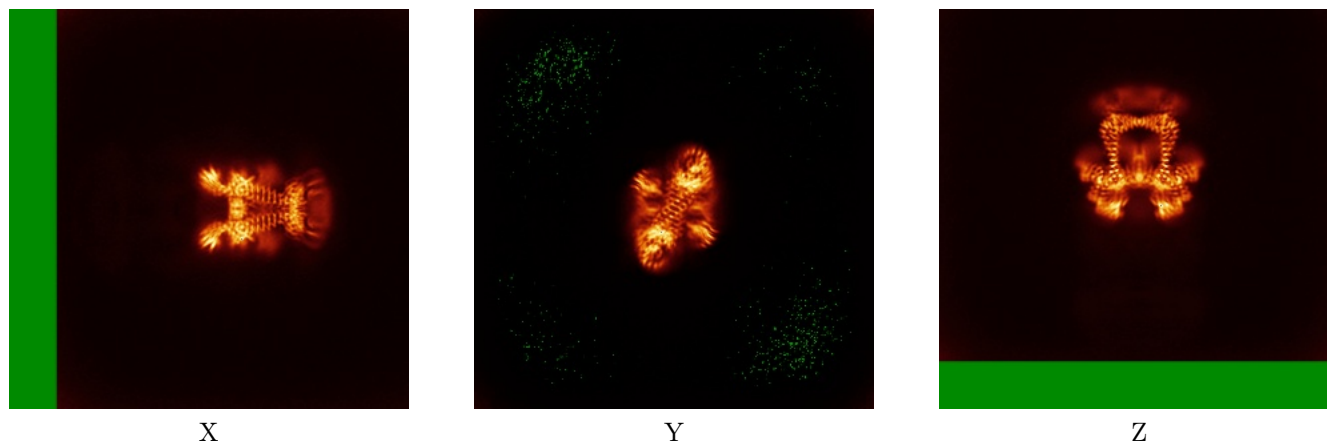


Z Index: 263

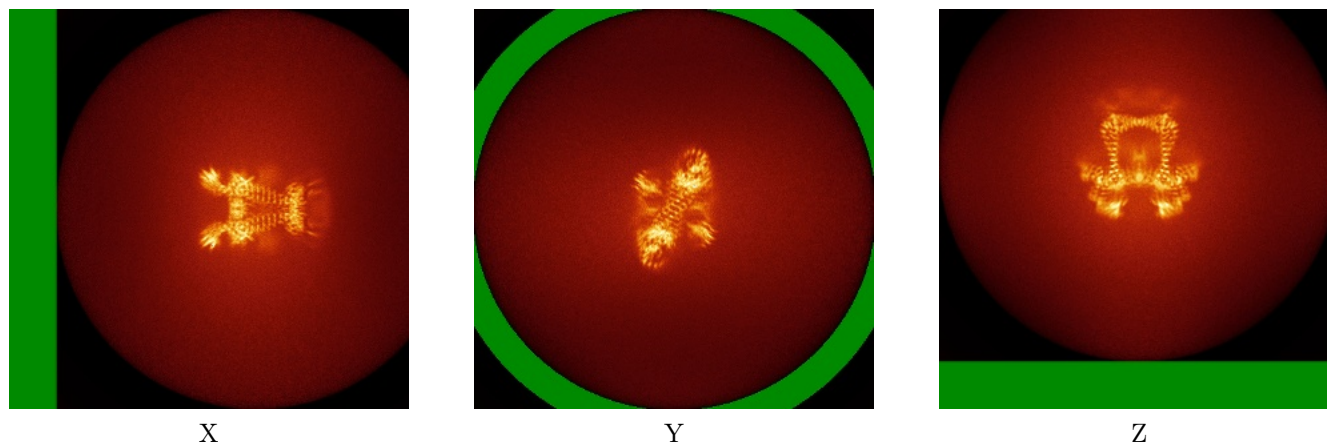
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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.

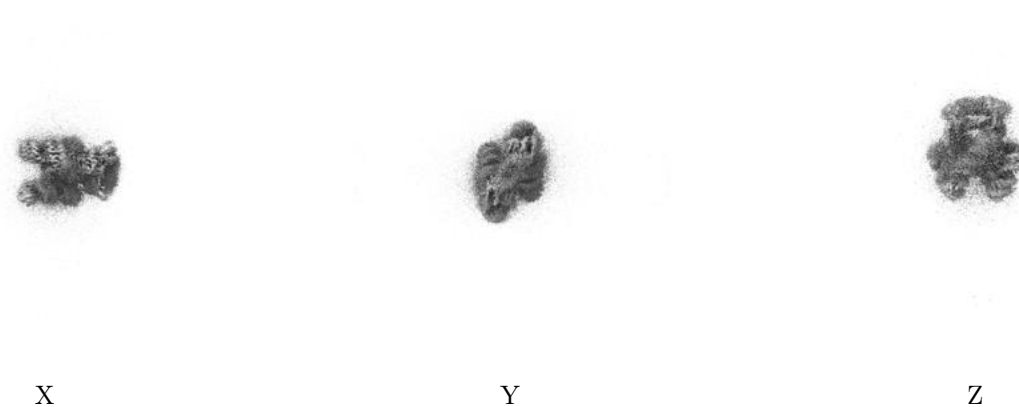
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.00115. 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.

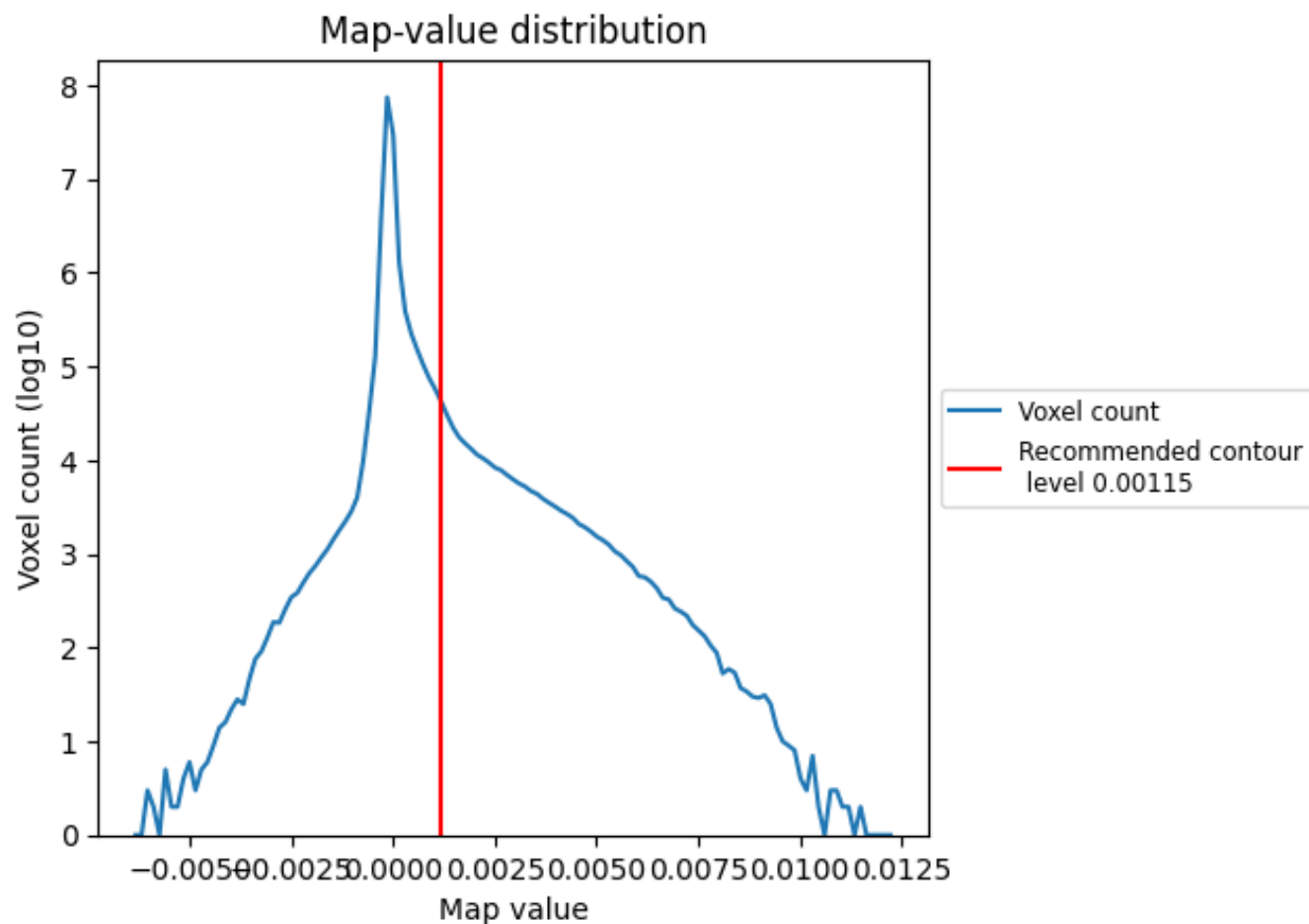
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

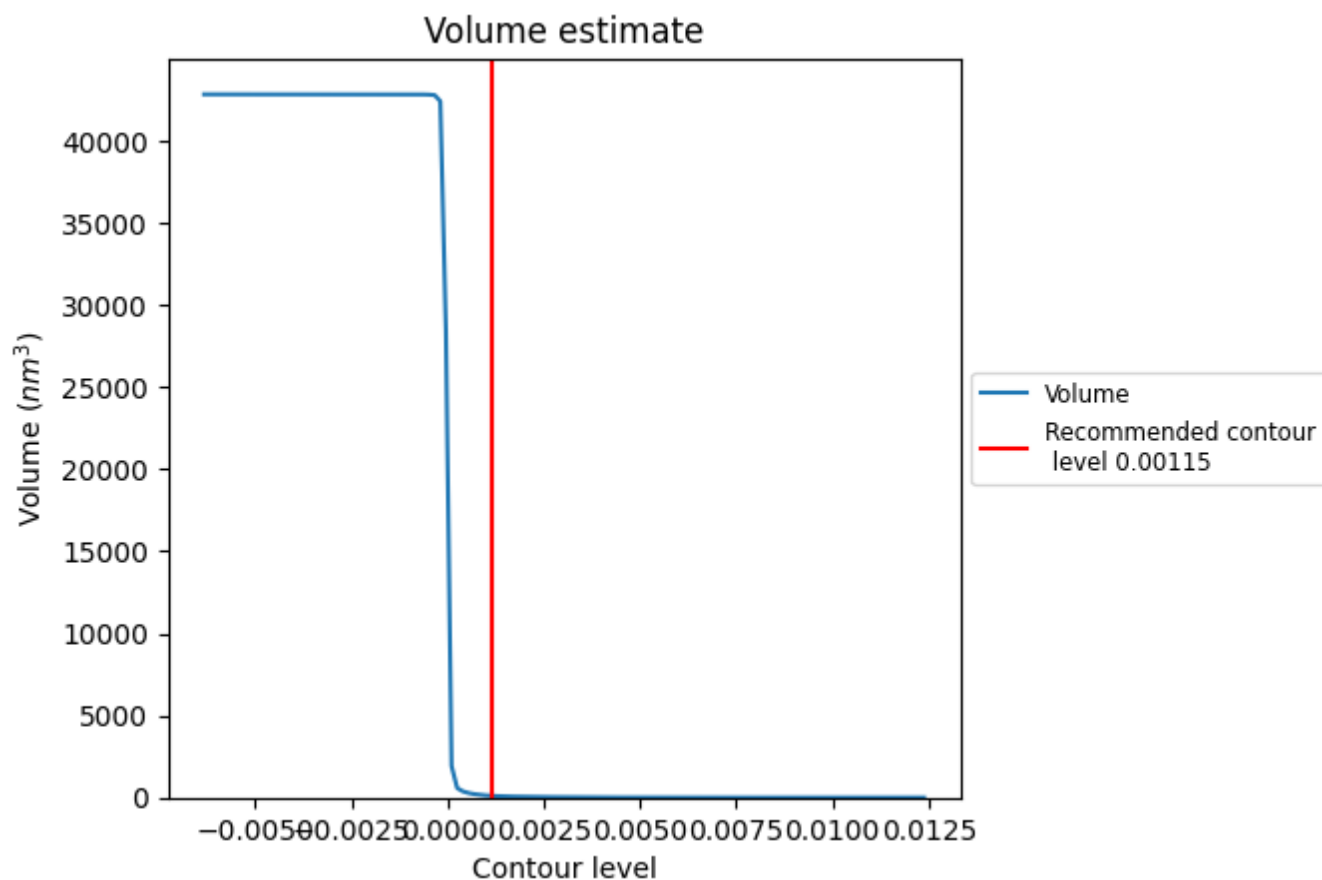
This section contains the results of statistical analysis of the map.

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.

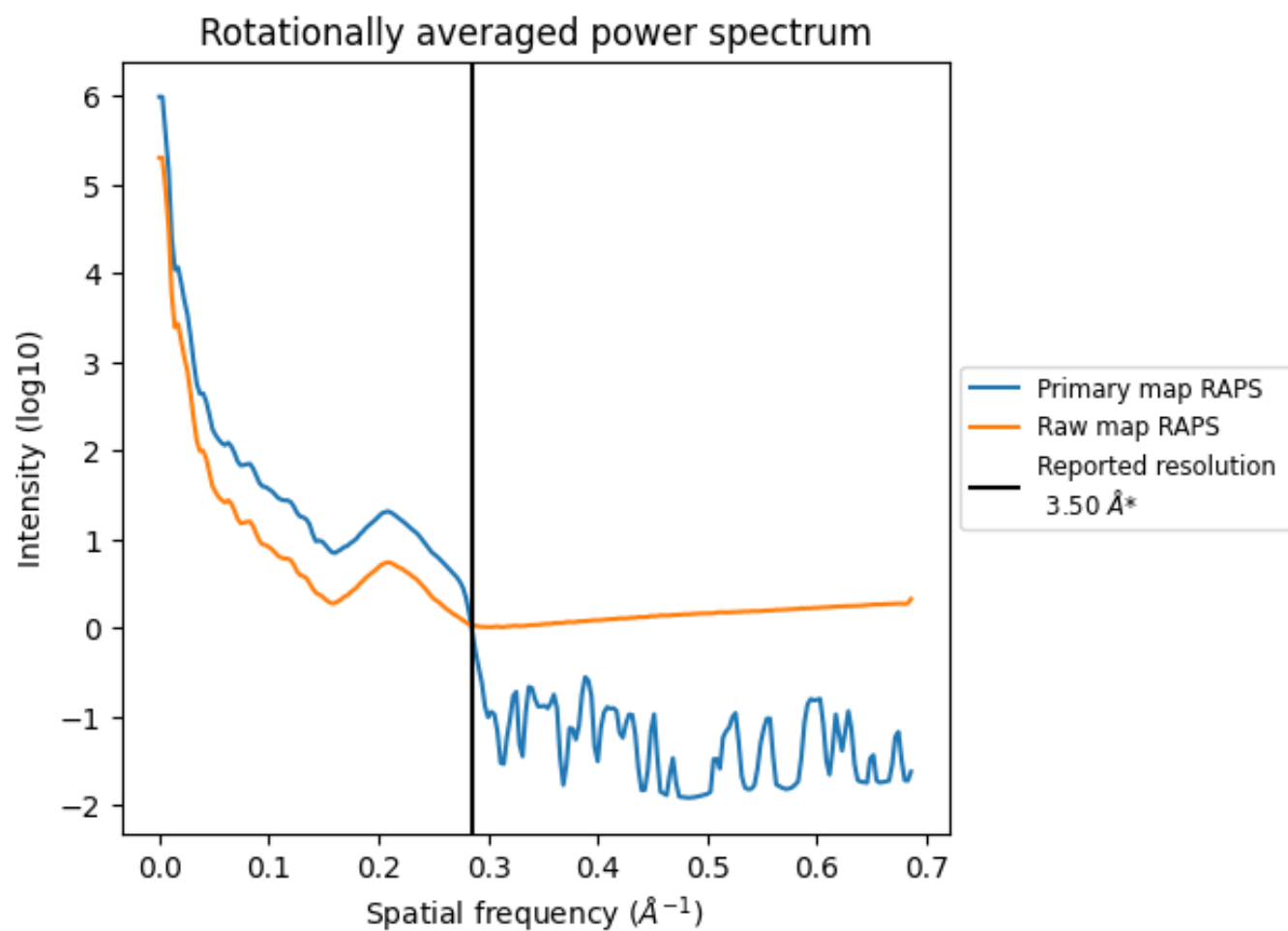
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 104 nm^3 ; this corresponds to an approximate mass of 94 kDa.

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

7.3 Rotationally averaged power spectrum [i](#)

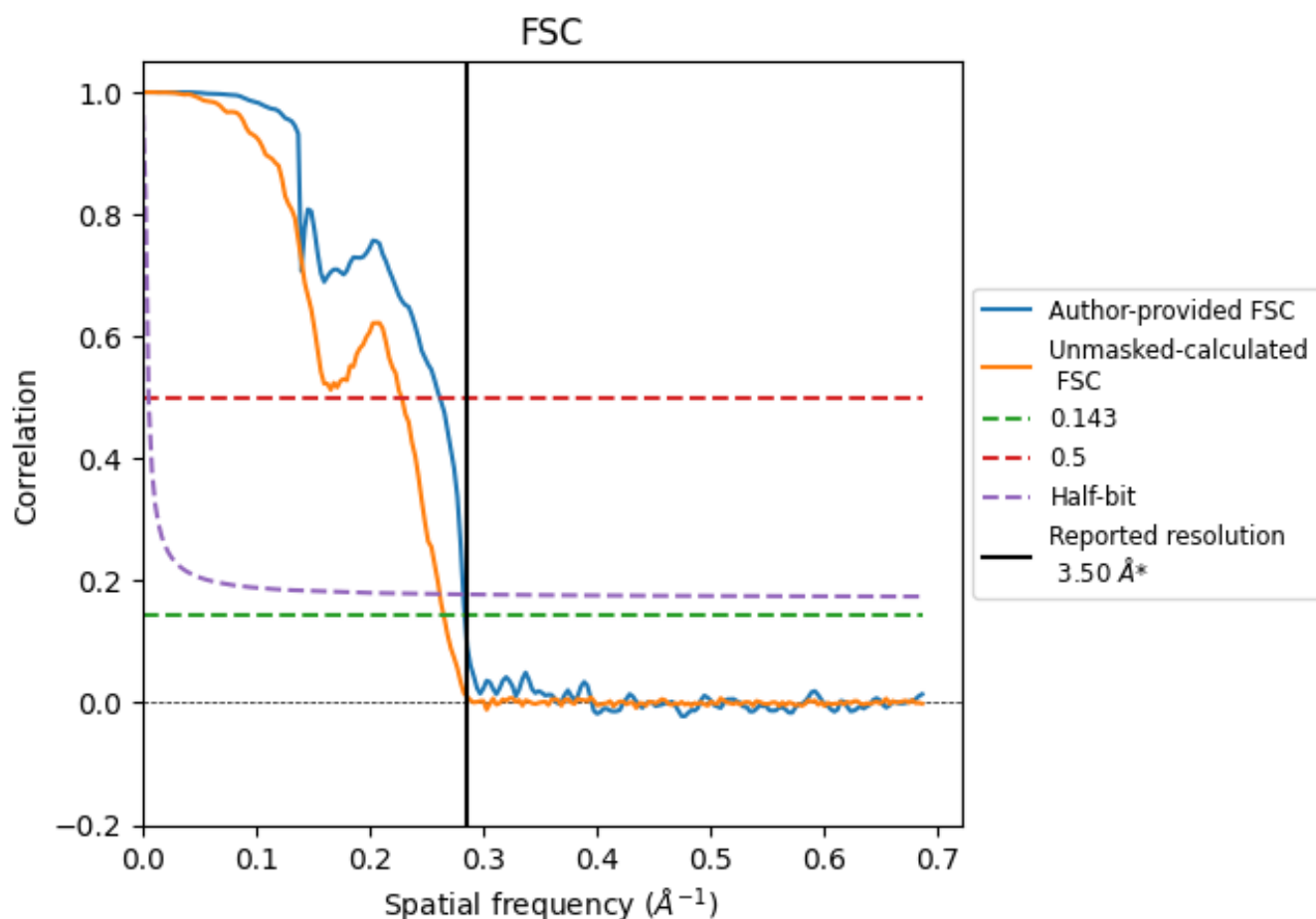


*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.286 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

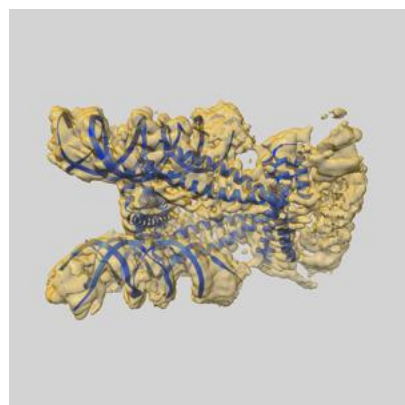
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	3.53	3.82	3.54
Unmasked-calculated*	3.77	4.38	3.82

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

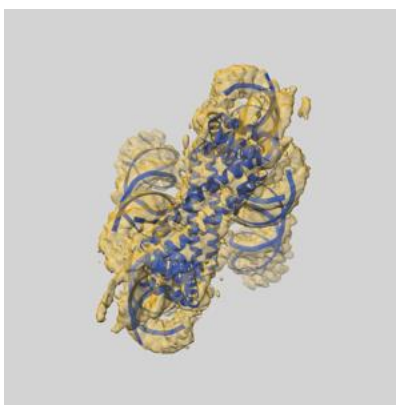
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-71807 and PDB model 9PRC. Per-residue inclusion information can be found in section [3](#) on page [5](#).

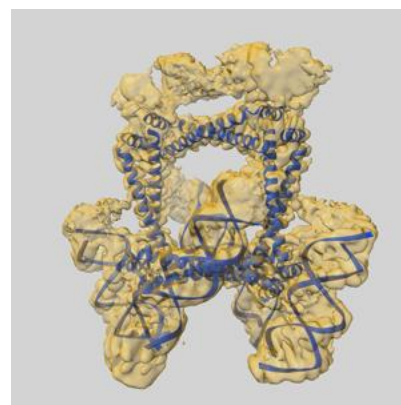
9.1 Map-model overlay [i](#)



X



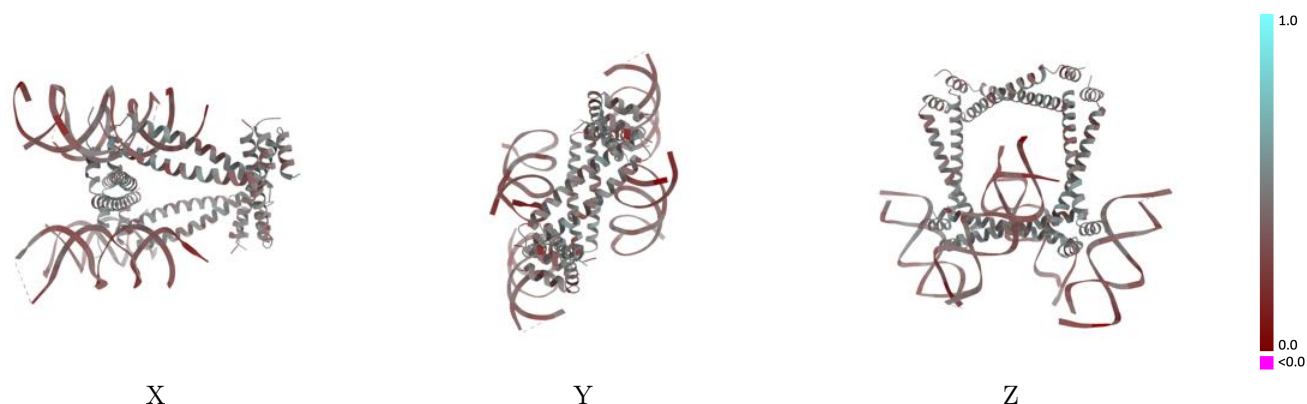
Y



Z

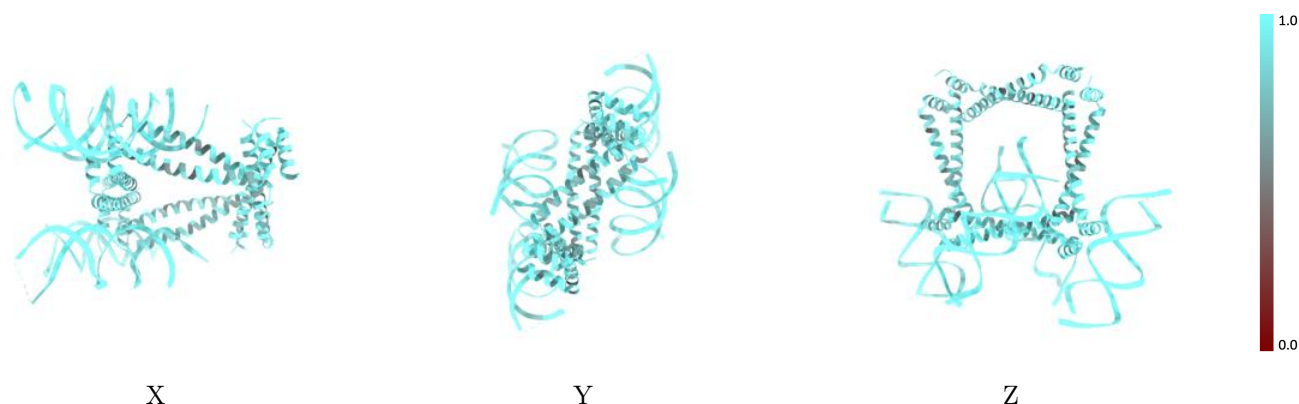
The images above show the 3D surface view of the map at the recommended contour level 0.00115 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



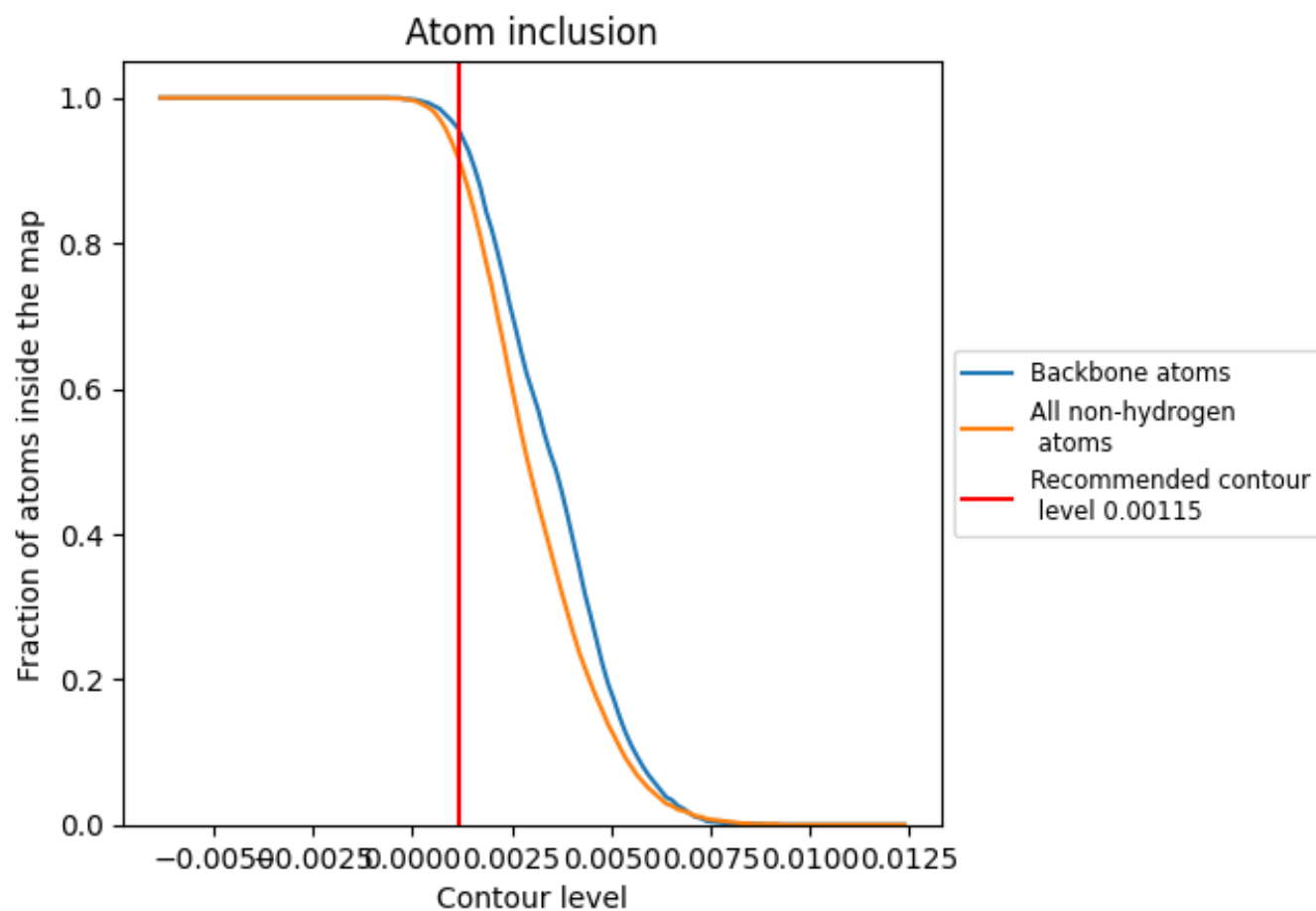
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.00115).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 92% 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 (0.00115) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9180	0.3920
A	0.9270	0.4680
B	0.9200	0.4700
C	0.8830	0.4550
D	0.8820	0.4620
E	0.8610	0.4370
F	0.8670	0.4430
G	0.8950	0.4700
H	0.8620	0.4410
J	0.9430	0.3350
K	0.9690	0.3210
M	0.9460	0.3470
P	0.9380	0.3050

