



Full wwPDB EM Validation Report ⓘ

Mar 24, 2026 – 09:59 PM UTC

PDB ID : 8HMZ / pdb_00008hmz
EMDB ID : EMD-34905
Title : Cryo-EM structure of the human post-catalytic TSEN/pre-tRNA complex
Authors : Zhang, X.; Yang, F.; Zhan, X.; Shi, Y.
Deposited on : 2022-12-06
Resolution : 2.90 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

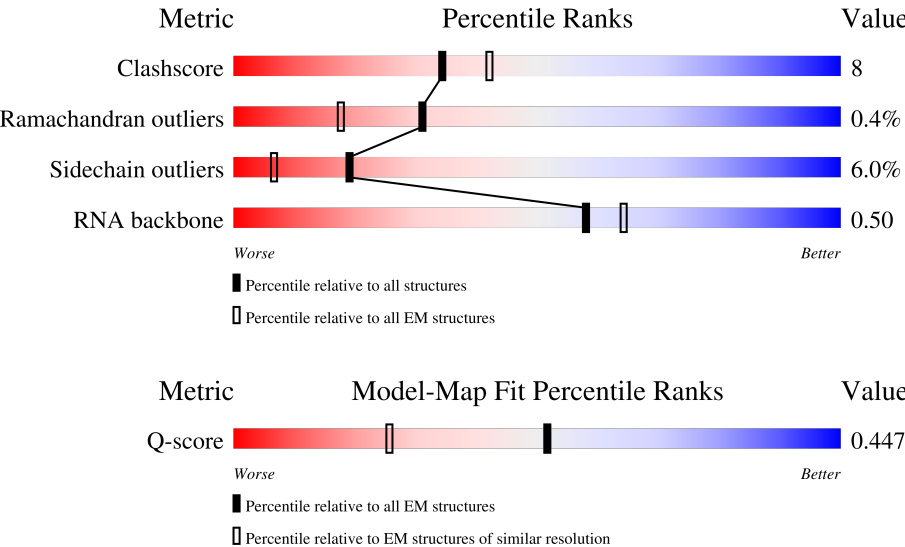
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13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

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.90 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	13054 (2.40 - 3.40)

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	485	25% 47% 16% 35%
2	B	330	5% 61% 18% 20%
3	C	546	41% 9% 48%

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Mol	Chain	Length	Quality of chain
4	D	213	42%16%41%
5	5	72	6%33%14%12%38%
6	3	42	71%14%12%
7	E	445	85%77%8%15%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 11542 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 tRNA-splicing endonuclease subunit Sen2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	313	Total	C	N	O	S	0	0
			2602	1680	443	466	13		

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q8NCE0
A	-18	ALA	-	expression tag	UNP Q8NCE0
A	-17	SER	-	expression tag	UNP Q8NCE0
A	-16	ASP	-	expression tag	UNP Q8NCE0
A	-15	TYR	-	expression tag	UNP Q8NCE0
A	-14	LYS	-	expression tag	UNP Q8NCE0
A	-13	ASP	-	expression tag	UNP Q8NCE0
A	-12	ASP	-	expression tag	UNP Q8NCE0
A	-11	ASP	-	expression tag	UNP Q8NCE0
A	-10	ASP	-	expression tag	UNP Q8NCE0
A	-9	LYS	-	expression tag	UNP Q8NCE0
A	-8	ALA	-	expression tag	UNP Q8NCE0
A	-7	SER	-	expression tag	UNP Q8NCE0
A	-6	ASP	-	expression tag	UNP Q8NCE0
A	-5	GLU	-	expression tag	UNP Q8NCE0
A	-4	VAL	-	expression tag	UNP Q8NCE0
A	-3	ASP	-	expression tag	UNP Q8NCE0
A	-2	ALA	-	expression tag	UNP Q8NCE0
A	-1	GLY	-	expression tag	UNP Q8NCE0
A	0	THR	-	expression tag	UNP Q8NCE0

- Molecule 2 is a protein called tRNA-splicing endonuclease subunit Sen34.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	264	Total	C	N	O	S	0	0
			2062	1301	388	369	4		

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP Q9BSV6
B	-18	ALA	-	expression tag	UNP Q9BSV6
B	-17	SER	-	expression tag	UNP Q9BSV6
B	-16	ASP	-	expression tag	UNP Q9BSV6
B	-15	TYR	-	expression tag	UNP Q9BSV6
B	-14	LYS	-	expression tag	UNP Q9BSV6
B	-13	ASP	-	expression tag	UNP Q9BSV6
B	-12	ASP	-	expression tag	UNP Q9BSV6
B	-11	ASP	-	expression tag	UNP Q9BSV6
B	-10	ASP	-	expression tag	UNP Q9BSV6
B	-9	LYS	-	expression tag	UNP Q9BSV6
B	-8	ALA	-	expression tag	UNP Q9BSV6
B	-7	SER	-	expression tag	UNP Q9BSV6
B	-6	ASP	-	expression tag	UNP Q9BSV6
B	-5	GLU	-	expression tag	UNP Q9BSV6
B	-4	VAL	-	expression tag	UNP Q9BSV6
B	-3	ASP	-	expression tag	UNP Q9BSV6
B	-2	ALA	-	expression tag	UNP Q9BSV6
B	-1	GLY	-	expression tag	UNP Q9BSV6
B	0	THR	-	expression tag	UNP Q9BSV6

- Molecule 3 is a protein called tRNA-splicing endonuclease subunit Sen54.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	285	Total	C	N	O	S	0	0
			2263	1440	405	412	6		

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-19	MET	-	initiating methionine	UNP Q7Z6J9
C	-18	ALA	-	expression tag	UNP Q7Z6J9
C	-17	SER	-	expression tag	UNP Q7Z6J9
C	-16	ASP	-	expression tag	UNP Q7Z6J9
C	-15	TYR	-	expression tag	UNP Q7Z6J9
C	-14	LYS	-	expression tag	UNP Q7Z6J9
C	-13	ASP	-	expression tag	UNP Q7Z6J9
C	-12	ASP	-	expression tag	UNP Q7Z6J9
C	-11	ASP	-	expression tag	UNP Q7Z6J9
C	-10	ASP	-	expression tag	UNP Q7Z6J9
C	-9	LYS	-	expression tag	UNP Q7Z6J9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	ALA	-	expression tag	UNP Q7Z6J9
C	-7	SER	-	expression tag	UNP Q7Z6J9
C	-6	ASP	-	expression tag	UNP Q7Z6J9
C	-5	GLU	-	expression tag	UNP Q7Z6J9
C	-4	VAL	-	expression tag	UNP Q7Z6J9
C	-3	ASP	-	expression tag	UNP Q7Z6J9
C	-2	ALA	-	expression tag	UNP Q7Z6J9
C	-1	GLY	-	expression tag	UNP Q7Z6J9
C	0	THR	-	expression tag	UNP Q7Z6J9

- Molecule 4 is a protein called Chromosome 1 open reading frame 19, isoform CRA_a.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	126	Total	C	N	O	S	0	0
			984	631	154	191	8		

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-37	MET	-	initiating methionine	UNP A0A2U3TZM3
D	-36	ALA	-	expression tag	UNP A0A2U3TZM3
D	-35	SER	-	expression tag	UNP A0A2U3TZM3
D	-34	SER	-	expression tag	UNP A0A2U3TZM3
D	-33	ALA	-	expression tag	UNP A0A2U3TZM3
D	-32	TRP	-	expression tag	UNP A0A2U3TZM3
D	-31	SER	-	expression tag	UNP A0A2U3TZM3
D	-30	HIS	-	expression tag	UNP A0A2U3TZM3
D	-29	PRO	-	expression tag	UNP A0A2U3TZM3
D	-28	GLN	-	expression tag	UNP A0A2U3TZM3
D	-27	PHE	-	expression tag	UNP A0A2U3TZM3
D	-26	GLU	-	expression tag	UNP A0A2U3TZM3
D	-25	LYS	-	expression tag	UNP A0A2U3TZM3
D	-24	GLY	-	expression tag	UNP A0A2U3TZM3
D	-23	GLY	-	expression tag	UNP A0A2U3TZM3
D	-22	GLY	-	expression tag	UNP A0A2U3TZM3
D	-21	SER	-	expression tag	UNP A0A2U3TZM3
D	-20	GLY	-	expression tag	UNP A0A2U3TZM3
D	-19	GLY	-	expression tag	UNP A0A2U3TZM3
D	-18	GLY	-	expression tag	UNP A0A2U3TZM3
D	-17	SER	-	expression tag	UNP A0A2U3TZM3
D	-16	GLY	-	expression tag	UNP A0A2U3TZM3
D	-15	GLY	-	expression tag	UNP A0A2U3TZM3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-14	SER	-	expression tag	UNP A0A2U3TZM3
D	-13	ALA	-	expression tag	UNP A0A2U3TZM3
D	-12	TRP	-	expression tag	UNP A0A2U3TZM3
D	-11	SER	-	expression tag	UNP A0A2U3TZM3
D	-10	HIS	-	expression tag	UNP A0A2U3TZM3
D	-9	PRO	-	expression tag	UNP A0A2U3TZM3
D	-8	GLN	-	expression tag	UNP A0A2U3TZM3
D	-7	PHE	-	expression tag	UNP A0A2U3TZM3
D	-6	GLU	-	expression tag	UNP A0A2U3TZM3
D	-5	LYS	-	expression tag	UNP A0A2U3TZM3
D	-4	GLY	-	expression tag	UNP A0A2U3TZM3
D	-3	SER	-	expression tag	UNP A0A2U3TZM3
D	-2	ALA	-	expression tag	UNP A0A2U3TZM3
D	-1	ALA	-	expression tag	UNP A0A2U3TZM3
D	0	ALA	-	expression tag	UNP A0A2U3TZM3

- Molecule 5 is a RNA chain called Pre-tRNA 5' END.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	5	45	Total	C	N	O	P	0	0
			968	430	174	318	46		

- Molecule 6 is a RNA chain called Pre-tRNA 3' END.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	3	37	Total	C	N	O	P	0	0
			788	353	142	257	36		

- Molecule 7 is a protein called Polyribonucleotide 5'-hydroxyl-kinase Clp1.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	E	379	Total	C	N	O	0	0
			1869	1111	379	379		

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-19	MET	-	initiating methionine	UNP Q92989
E	-18	ALA	-	expression tag	UNP Q92989
E	-17	SER	-	expression tag	UNP Q92989
E	-16	ASP	-	expression tag	UNP Q92989
E	-15	TYR	-	expression tag	UNP Q92989

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-14	LYS	-	expression tag	UNP Q92989
E	-13	ASP	-	expression tag	UNP Q92989
E	-12	ASP	-	expression tag	UNP Q92989
E	-11	ASP	-	expression tag	UNP Q92989
E	-10	ASP	-	expression tag	UNP Q92989
E	-9	LYS	-	expression tag	UNP Q92989
E	-8	ALA	-	expression tag	UNP Q92989
E	-7	SER	-	expression tag	UNP Q92989
E	-6	ASP	-	expression tag	UNP Q92989
E	-5	GLU	-	expression tag	UNP Q92989
E	-4	VAL	-	expression tag	UNP Q92989
E	-3	ASP	-	expression tag	UNP Q92989
E	-2	ALA	-	expression tag	UNP Q92989
E	-1	GLY	-	expression tag	UNP Q92989
E	0	THR	-	expression tag	UNP Q92989

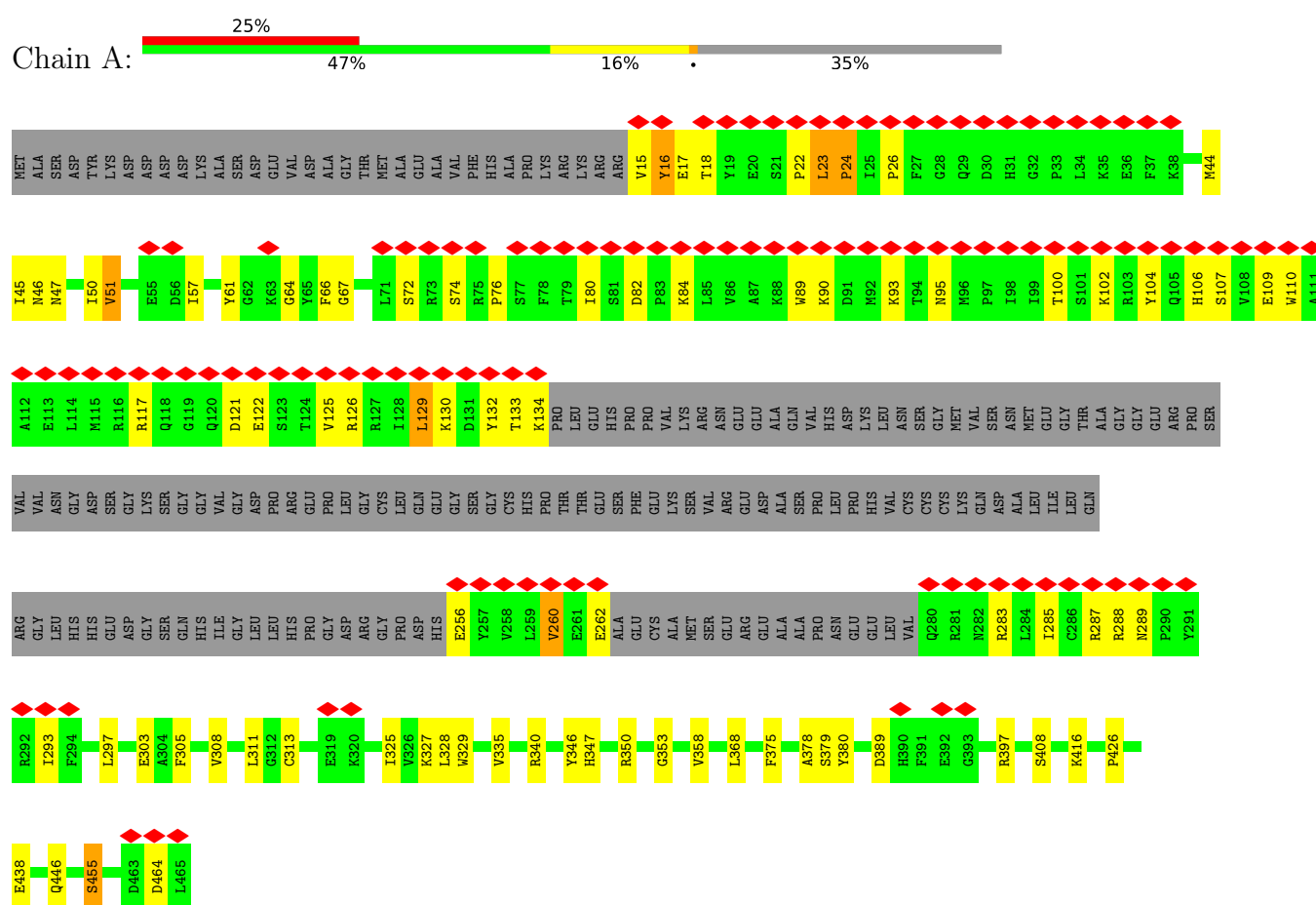
- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	AltConf
8	5	3	Total Mg 3 3	0
8	3	3	Total Mg 3 3	0

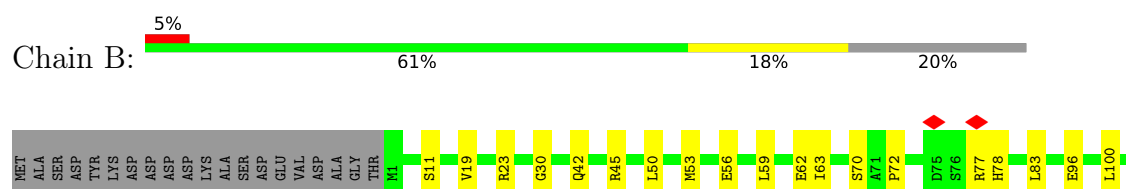
3 Residue-property plots

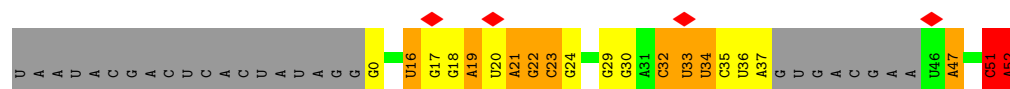
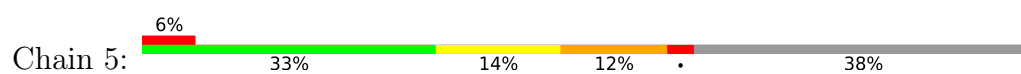
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: tRNA-splicing endonuclease subunit Sen2

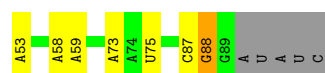


• Molecule 2: tRNA-splicing endonuclease subunit Sen34

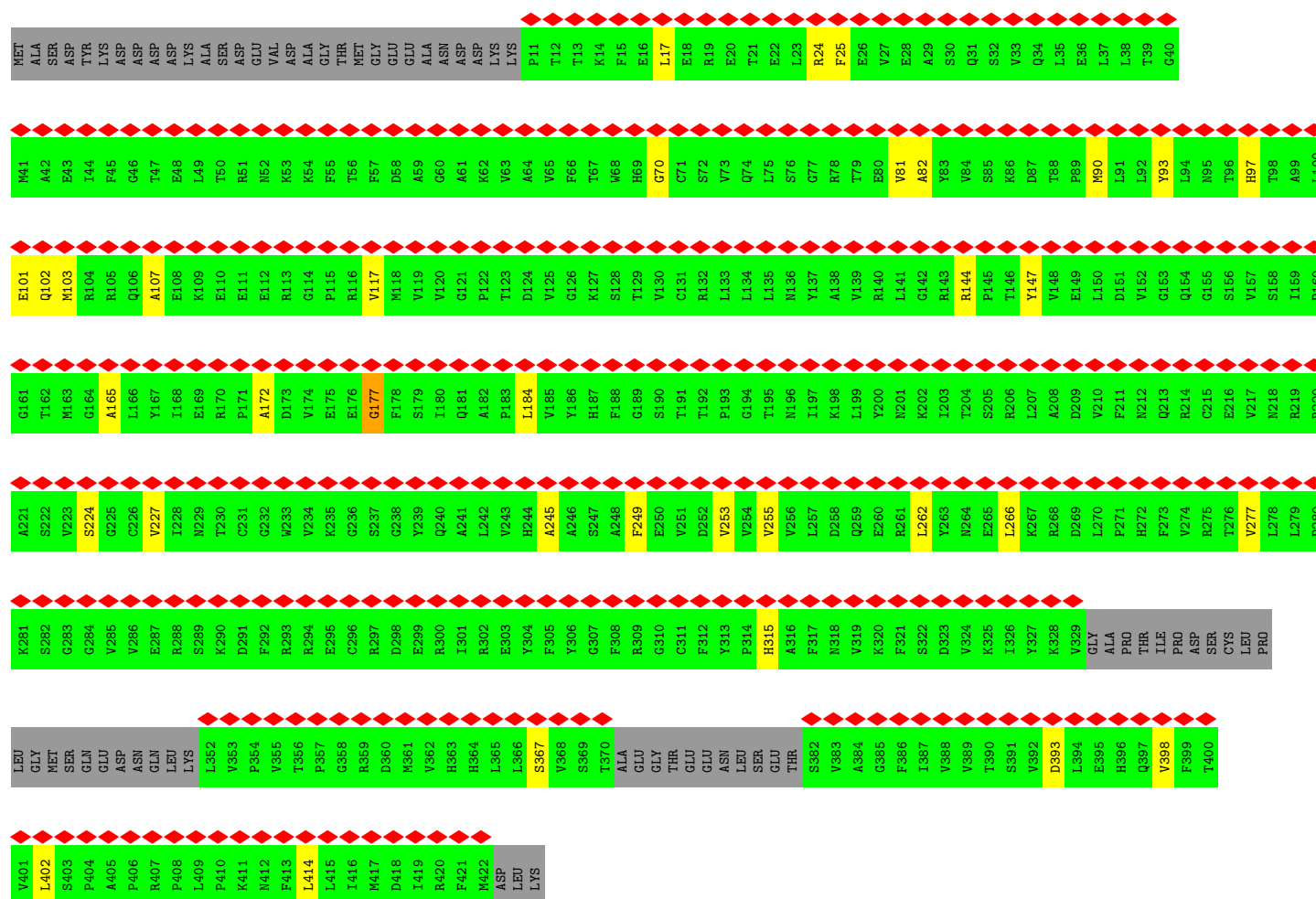
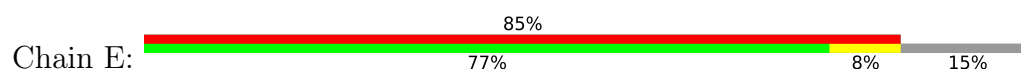




• Molecule 6: Pre-tRNA 3' END



• Molecule 7: Polyrribonucleotide 5'-hydroxyl-kinase Clp1



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	309195	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50.00	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	8.282	Depositor
Minimum map value	-4.794	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.133	Depositor
Recommended contour level	0.6	Depositor
Map size ($\text{\AA}$)	304.36002, 304.36002, 304.36002	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.087, 1.087, 1.087	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, A23

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	0/2665	0.47	0/3596
2	B	0.18	0/2108	0.43	0/2856
3	C	0.37	0/2319	0.85	23/3140 (0.7%)
4	D	0.17	0/1004	0.42	0/1366
5	5	0.41	2/1053 (0.2%)	0.43	2/1638 (0.1%)
6	3	0.14	0/881	0.23	0/1372
7	E	0.24	0/1866	0.65	1/2592 (0.0%)
All	All	0.26	2/11896 (0.0%)	0.57	26/16560 (0.2%)

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
3	C	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	5	51	C	P-OP2	-7.96	1.33	1.49
5	5	51	C	P-OP1	-5.85	1.37	1.49

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	412	GLY	CA-C-N	10.44	141.48	121.54
3	C	412	GLY	C-N-CA	10.44	141.48	121.54
5	5	51	C	O4'-C1'-C2'	-9.37	96.43	105.80
3	C	414	ALA	O-C-N	8.34	133.69	122.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	420	VAL	CA-C-N	8.32	137.43	121.54
3	C	420	VAL	C-N-CA	8.32	137.43	121.54
3	C	419	VAL	CA-C-N	8.17	136.67	121.97
3	C	419	VAL	C-N-CA	8.17	136.67	121.97
3	C	421	LEU	CA-C-N	-7.84	110.10	122.73
3	C	421	LEU	C-N-CA	-7.84	110.10	122.73
3	C	414	ALA	CA-C-N	7.58	136.03	121.54
3	C	414	ALA	C-N-CA	7.58	136.03	121.54
3	C	419	VAL	N-CA-CB	7.29	120.64	111.46
3	C	521	PRO	O-C-N	-6.85	117.87	122.73
3	C	523	ASP	CA-C-O	-6.71	112.68	120.20
5	5	51	C	C4'-C3'-O3'	6.42	119.03	109.40
3	C	418	ALA	CA-C-N	-6.28	114.96	123.12
3	C	418	ALA	C-N-CA	-6.28	114.96	123.12
3	C	420	VAL	O-C-N	6.11	130.20	122.57
3	C	411	PRO	CA-N-CD	-5.72	103.99	112.00
3	C	420	VAL	CA-C-O	-5.71	113.65	120.78
3	C	413	GLN	CB-CA-C	-5.70	99.08	110.42
3	C	420	VAL	N-CA-C	-5.68	97.53	109.34
3	C	523	ASP	O-C-N	5.55	128.53	122.20
3	C	417	PRO	CA-C-O	-5.41	117.31	121.31
7	E	177	GLY	N-CA-C	5.15	118.12	110.60

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	412	GLY	Mainchain
3	C	414	ALA	Mainchain
3	C	521	PRO	Mainchain
3	C	522	GLN	Mainchain

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	2602	0	2626	54	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2062	0	2091	34	0
3	C	2263	0	2234	44	0
4	D	984	0	987	23	0
5	5	968	0	484	17	0
6	3	788	0	400	7	0
7	E	1869	0	827	20	0
8	3	3	0	0	0	0
8	5	3	0	0	0	0
All	All	11542	0	9649	171	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 8.

All (171) 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)
7:E:17:LEU:O	7:E:70:GLY:HA2	1.77	0.84
5:5:36:U:O2	5:5:47:A:N6	2.10	0.84
1:A:311:LEU:HD21	3:C:414:ALA:HB2	1.61	0.81
1:A:23:LEU:HB2	1:A:26:PRO:HD2	1.67	0.77
7:E:25:PHE:HA	7:E:81:VAL:O	1.85	0.76
2:B:248:PRO:HB2	3:C:45:PRO:HG3	1.68	0.75
7:E:367:SER:O	7:E:414:LEU:HA	1.88	0.74
1:A:346:TYR:OH	1:A:350:ARG:NH1	2.23	0.72
1:A:378:ALA:HB3	1:A:416:LYS:HE2	1.72	0.70
1:A:44:MET:HE3	1:A:47:ASN:HA	1.74	0.69
2:B:53:MET:HB2	2:B:56:GLU:HG3	1.75	0.69
2:B:62:GLU:OE1	2:B:221:TYR:OH	2.10	0.69
5:5:52:A23:O1C	6:3:53:A:O5'	2.09	0.69
7:E:315:HIS:O	7:E:402:LEU:HA	1.94	0.68
2:B:237:ALA:HB2	2:B:245:LEU:HG	1.74	0.68
3:C:440:LEU:O	3:C:443:SER:OG	2.12	0.67
5:5:16:U:O2'	5:5:19:A:OP1	2.11	0.67
4:D:156:THR:HG21	4:D:160:MET:HE2	1.75	0.66
4:D:49:MET:HG2	4:D:90:ILE:HD11	1.78	0.66
2:B:205:GLN:NE2	2:B:212:ALA:O	2.29	0.65
3:C:413:GLN:H	3:C:413:GLN:CD	2.04	0.65
1:A:16:TYR:O	1:A:18:THR:N	2.30	0.65
7:E:172:ALA:HB1	7:E:177:GLY:HA2	1.79	0.65
4:D:138:SER:HB3	4:D:156:THR:HG23	1.78	0.64
2:B:250:ASP:OD2	3:C:59:ARG:NH2	2.30	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:5:52:A23:PC	6:3:53:A:O5'	2.56	0.63
7:E:97:HIS:O	7:E:101:GLU:N	2.30	0.63
1:A:260:VAL:HB	1:A:287:ARG:HG3	1.81	0.62
2:B:23:ARG:NH1	2:B:250:ASP:OD1	2.32	0.62
4:D:120:ARG:NH2	4:D:157:ASP:OD1	2.33	0.62
3:C:413:GLN:CD	3:C:413:GLN:N	2.58	0.62
3:C:421:LEU:HD22	3:C:421:LEU:O	1.99	0.61
1:A:389:ASP:HB2	1:A:426:PRO:HA	1.81	0.61
1:A:57:ILE:HG12	1:A:297:LEU:HB2	1.82	0.61
2:B:184:LEU:HD21	3:C:27:ARG:HD2	1.82	0.61
3:C:167:VAL:HB	3:C:453:TYR:HB2	1.84	0.60
2:B:206:SER:OG	2:B:207:LYS:N	2.34	0.60
3:C:472:ARG:NH1	3:C:495:SER:O	2.36	0.58
1:A:358:VAL:HG21	4:D:145:GLU:HG2	1.85	0.58
3:C:73:ARG:NH2	5:5:16:U:O5'	2.37	0.58
1:A:64:GLY:O	1:A:350:ARG:NH2	2.37	0.57
1:A:368:LEU:HD12	1:A:380:TYR:HB2	1.87	0.57
1:A:15:VAL:HB	1:A:24:PRO:HA	1.85	0.57
6:3:87:C:H2'	6:3:88:G:O4'	2.05	0.57
2:B:256:ALA:HB3	2:B:286:LYS:HE3	1.88	0.56
2:B:83:LEU:HD11	3:C:132:ASP:HB3	1.88	0.56
7:E:147:TYR:O	7:E:165:ALA:HA	2.06	0.56
2:B:184:LEU:HD22	3:C:71:VAL:HG12	1.89	0.55
7:E:90:MET:O	7:E:93:TYR:CB	2.54	0.55
7:E:24:ARG:O	7:E:82:ALA:HA	2.06	0.55
1:A:47:ASN:HB3	1:A:335:VAL:HG21	1.89	0.55
6:3:73:A:O2'	6:3:75:U:OP2	2.22	0.55
1:A:106:HIS:HB3	1:A:110:TRP:CH2	2.42	0.55
4:D:116:HIS:O	4:D:120:ARG:HG2	2.07	0.54
3:C:99:TRP:CH2	3:C:110:GLN:HB3	2.42	0.54
2:B:77:ARG:HG3	2:B:78:HIS:CD2	2.42	0.54
2:B:198:ARG:HG3	2:B:199:PRO:HD2	1.89	0.53
3:C:476:SER:OG	3:C:477:GLY:N	2.41	0.53
7:E:117:VAL:HA	7:E:253:VAL:O	2.09	0.53
1:A:129:LEU:O	1:A:133:THR:HG23	2.09	0.53
2:B:130:ALA:O	2:B:134:SER:OG	2.23	0.53
6:3:58:A:H2'	6:3:59:A:C8	2.44	0.53
1:A:256:GLU:O	1:A:289:ASN:ND2	2.42	0.52
4:D:66:VAL:HG21	4:D:107:PRO:HB3	1.91	0.52
3:C:421:LEU:O	3:C:421:LEU:HD13	2.10	0.52
3:C:521:PRO:C	3:C:523:ASP:H	2.17	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:413:GLN:H	3:C:413:GLN:NE2	2.08	0.51
3:C:122:GLU:OE2	3:C:155:TYR:OH	2.23	0.51
7:E:255:VAL:HA	7:E:277:VAL:O	2.10	0.51
2:B:126:LYS:O	2:B:129:GLN:HG3	2.10	0.51
1:A:106:HIS:HA	1:A:109:GLU:HG2	1.91	0.51
1:A:67:GLY:N	1:A:303:GLU:OE1	2.43	0.51
1:A:107:SER:HA	1:A:110:TRP:NE1	2.26	0.51
4:D:145:GLU:HG3	4:D:146:SER:H	1.76	0.51
3:C:32:LYS:HD3	5:5:23:C:H5'	1.92	0.51
1:A:15:VAL:HG22	1:A:325:ILE:HG23	1.94	0.50
1:A:80:ILE:HG21	1:A:293:ILE:HD13	1.93	0.50
1:A:262:GLU:N	1:A:283:ARG:O	2.45	0.50
5:5:0:G:H1	6:3:88:G:H21	1.58	0.50
1:A:46:ASN:ND2	4:D:46:TYR:OH	2.44	0.50
1:A:311:LEU:HD22	1:A:313:CYS:SG	2.51	0.50
1:A:455:SER:O	3:C:166:TYR:OH	2.29	0.50
3:C:522:GLN:O	3:C:522:GLN:HG3	2.07	0.49
1:A:76:PRO:HG3	1:A:297:LEU:HD13	1.94	0.49
3:C:30:SER:OG	3:C:31:GLN:N	2.45	0.49
2:B:287:THR:HG23	2:B:305:GLN:HG2	1.95	0.49
7:E:245:ALA:O	7:E:249:PHE:N	2.37	0.49
4:D:49:MET:HB3	4:D:60:VAL:HG23	1.94	0.48
5:5:51:C:H2'	5:5:51:C:O2	2.13	0.47
1:A:325:ILE:HA	1:A:328:LEU:HB2	1.97	0.47
4:D:135:LEU:HD23	4:D:135:LEU:H	1.78	0.47
1:A:397:ARG:NH1	2:B:272:GLN:OE1	2.47	0.47
1:A:104:TYR:HA	1:A:107:SER:OG	2.14	0.47
7:E:144:ARG:H	7:E:224:SER:CB	2.28	0.47
7:E:393:ASP:CB	7:E:398:VAL:O	2.63	0.47
1:A:15:VAL:O	1:A:15:VAL:HG12	2.15	0.46
2:B:123:LYS:O	2:B:126:LYS:HG3	2.15	0.46
1:A:329:TRP:HZ2	1:A:340:ARG:HD3	1.81	0.46
1:A:353:GLY:HA3	3:C:524:VAL:HG23	1.97	0.46
2:B:292:SER:HB3	4:D:116:HIS:ND1	2.31	0.46
5:5:32:C:H4'	5:5:33:U:OP2	2.16	0.46
2:B:304:LEU:HD12	4:D:155:LEU:HG	1.96	0.46
1:A:107:SER:HA	1:A:110:TRP:CD1	2.50	0.46
2:B:191:ARG:O	3:C:14:ALA:HB1	2.15	0.46
3:C:430:THR:HG22	3:C:431:HIS:H	1.81	0.46
7:E:17:LEU:O	7:E:70:GLY:CA	2.58	0.46
3:C:421:LEU:HD22	3:C:421:LEU:C	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:5:34:U:H2'	5:5:35:C:C6	2.50	0.46
3:C:484:ASP:N	3:C:484:ASP:OD1	2.46	0.45
5:5:52:A23:PC	6:3:53:A:HO5'	2.39	0.45
3:C:446:LEU:HD13	3:C:502:PHE:HE1	1.81	0.45
1:A:66:PHE:CE2	3:C:413:GLN:NE2	2.85	0.45
2:B:219:LEU:O	2:B:223:ILE:HD12	2.17	0.45
1:A:16:TYR:HA	1:A:22:PRO:HA	1.98	0.45
7:E:262:LEU:O	7:E:266:LEU:CB	2.65	0.45
2:B:30:GLY:HA3	2:B:50:LEU:HD21	1.99	0.45
3:C:21:ARG:HH11	3:C:21:ARG:HG2	1.82	0.45
3:C:521:PRO:C	3:C:523:ASP:N	2.74	0.45
1:A:82:ASP:OD2	1:A:100:THR:HG23	2.17	0.44
2:B:59:LEU:O	2:B:63:ILE:HG23	2.18	0.44
3:C:414:ALA:C	3:C:416:SER:H	2.26	0.44
1:A:133:THR:HB	1:A:287:ARG:NH1	2.33	0.44
2:B:72:PRO:HG3	2:B:197:ALA:HB2	2.00	0.44
7:E:102:GLN:O	7:E:103:MET:C	2.61	0.44
1:A:305:PHE:HA	1:A:308:VAL:HG12	2.00	0.43
1:A:23:LEU:H	1:A:23:LEU:HD22	1.84	0.43
2:B:180:ARG:HD2	2:B:183:LEU:HD12	1.99	0.43
2:B:253:ARG:HG3	3:C:55:LEU:HD13	2.01	0.43
3:C:421:LEU:O	3:C:421:LEU:CG	2.66	0.43
7:E:103:MET:O	7:E:107:ALA:N	2.50	0.43
1:A:61:TYR:OH	1:A:74:SER:O	2.16	0.43
3:C:27:ARG:NH1	5:5:16:U:OP1	2.52	0.43
1:A:117:ARG:HH12	4:D:47:LEU:HB2	1.83	0.43
4:D:74:LYS:HE3	4:D:152:TYR:CE2	2.54	0.43
3:C:474:CYS:SG	3:C:500:LEU:HD11	2.59	0.43
1:A:347:HIS:CE1	3:C:424:ILE:HG22	2.54	0.43
2:B:301:TYR:O	4:D:120:ARG:NH1	2.52	0.43
4:D:126:SER:O	4:D:130:GLN:HG2	2.18	0.43
5:5:35:C:H2'	5:5:36:U:C6	2.53	0.43
4:D:91:CYS:SG	4:D:104:THR:HG23	2.58	0.43
1:A:45:ILE:HG22	1:A:50:ILE:HG12	2.01	0.42
5:5:21:A:H5'	5:5:22:G:OP1	2.17	0.42
4:D:105:VAL:HG22	4:D:140:THR:HB	2.00	0.42
3:C:32:LYS:HB2	5:5:23:C:OP1	2.19	0.42
3:C:85:GLU:HG2	3:C:86:GLU:N	2.34	0.42
1:A:16:TYR:CD1	1:A:16:TYR:N	2.87	0.42
4:D:144:VAL:HG13	4:D:150:ILE:HG12	2.01	0.42
4:D:123:LEU:HD13	4:D:157:ASP:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:LYS:NZ	1:A:288:ARG:HH21	2.17	0.42
2:B:289:LEU:HD21	4:D:159:PHE:CE1	2.55	0.42
2:B:19:VAL:HG13	2:B:50:LEU:HD13	2.01	0.42
4:D:76:TRP:C	4:D:77:HIS:HD2	2.28	0.42
3:C:432:LEU:HD12	3:C:432:LEU:HA	1.83	0.41
7:E:147:TYR:HA	7:E:227:VAL:O	2.19	0.41
3:C:28:SER:O	3:C:28:SER:OG	2.35	0.41
5:5:23:C:H2'	5:5:24:G:C8	2.56	0.41
1:A:51:VAL:HG13	1:A:297:LEU:HB3	2.03	0.41
1:A:121:ASP:O	1:A:125:VAL:HG23	2.21	0.41
3:C:421:LEU:O	3:C:421:LEU:CD2	2.66	0.41
7:E:165:ALA:HB3	7:E:184:LEU:CB	2.51	0.41
1:A:84:LYS:HD2	1:A:84:LYS:O	2.21	0.41
1:A:90:LYS:O	1:A:90:LYS:HG2	2.21	0.41
2:B:214:ARG:O	2:B:218:GLU:HG2	2.20	0.41
1:A:122:GLU:HB3	1:A:126:ARG:CZ	2.51	0.40
5:5:23:C:H2'	5:5:24:G:H8	1.86	0.40
1:A:126:ARG:O	1:A:130:LYS:HB2	2.20	0.40
1:A:129:LEU:O	1:A:133:THR:N	2.52	0.40
2:B:200:LEU:HD21	2:B:209:TRP:HB2	2.04	0.40
2:B:280:LEU:O	2:B:284:VAL:HG22	2.21	0.40
3:C:85:GLU:CD	3:C:85:GLU:H	2.29	0.40
1:A:93:LYS:HA	1:A:93:LYS:HD3	1.80	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	307/485 (63%)	276 (90%)	29 (9%)	2 (1%)	18 48
2	B	260/330 (79%)	250 (96%)	10 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	281/546 (52%)	259 (92%)	19 (7%)	3 (1%)	11	36
4	D	124/213 (58%)	118 (95%)	5 (4%)	1 (1%)	16	44
7	E	373/445 (84%)	346 (93%)	27 (7%)	0	100	100
All	All	1345/2019 (67%)	1249 (93%)	90 (7%)	6 (0%)	31	59

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	GLU
3	C	413	GLN
3	C	421	LEU
3	C	524	VAL
4	D	162	PRO
1	A	24	PRO

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	288/428 (67%)	269 (93%)	19 (7%)	15	43
2	B	212/261 (81%)	201 (95%)	11 (5%)	21	52
3	C	242/464 (52%)	227 (94%)	15 (6%)	16	46
4	D	113/171 (66%)	107 (95%)	6 (5%)	20	52
All	All	855/1324 (65%)	804 (94%)	51 (6%)	19	47

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	TYR
1	A	23	LEU
1	A	51	VAL
1	A	72	SER
1	A	89	TRP

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Mol	Chain	Res	Type
1	A	95	ASN
1	A	102	LYS
1	A	129	LEU
1	A	132	TYR
1	A	260	VAL
1	A	285	ILE
1	A	327	LYS
1	A	375	PHE
1	A	379	SER
1	A	408	SER
1	A	438	GLU
1	A	446	GLN
1	A	455	SER
1	A	464	ASP
2	B	11	SER
2	B	42	GLN
2	B	45	ARG
2	B	70	SER
2	B	96	GLU
2	B	100	LEU
2	B	115	LYS
2	B	126	LYS
2	B	185	VAL
2	B	234	LEU
2	B	268	THR
3	C	100	GLN
3	C	176	VAL
3	C	413	GLN
3	C	420	VAL
3	C	421	LEU
3	C	425	SER
3	C	429	THR
3	C	430	THR
3	C	438	ARG
3	C	463	LYS
3	C	484	ASP
3	C	510	ILE
3	C	511	SER
3	C	522	GLN
3	C	523	ASP
4	D	48	GLU
4	D	81	CYS

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Mol	Chain	Res	Type
4	D	91	CYS
4	D	97	ILE
4	D	111	THR
4	D	130	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	59	GLN
1	A	120	GLN
1	A	446	GLN
2	B	89	GLN
2	B	90	GLN
2	B	120	GLN
2	B	186	GLN
2	B	205	GLN
3	C	69	GLN
3	C	138	GLN
3	C	172	GLN
3	C	454	GLN
3	C	507	HIS
3	C	522	GLN
4	D	130	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
5	5	43/72 (59%)	17 (39%)	0
6	3	36/42 (85%)	1 (2%)	0
All	All	79/114 (69%)	18 (22%)	0

All (18) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	5	16	U
5	5	17	G
5	5	18	G
5	5	19	A
5	5	20	U

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Mol	Chain	Res	Type
5	5	21	A
5	5	22	G
5	5	23	C
5	5	29	G
5	5	30	G
5	5	32	C
5	5	33	U
5	5	34	U
5	5	37	A
5	5	47	A
5	5	51	C
5	5	52	A23
6	3	88	G

There are no RNA pucker outliers to report.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	A23	5	52	5	23,28,29	1.67	6 (26%)	32,43,46	2.27	13 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	A23	5	52	5	-	0/7/35/36	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	5	52	A23	C5-C4	4.74	1.47	1.39
5	5	52	A23	C5'-C4'	-2.90	1.42	1.51
5	5	52	A23	C5-C6	2.77	1.48	1.41
5	5	52	A23	C8-N7	2.38	1.36	1.31
5	5	52	A23	C5-N7	-2.22	1.35	1.39
5	5	52	A23	O5'-C5'	-2.18	1.37	1.44

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	5	52	A23	C5-C4-N3	-5.83	118.69	126.72
5	5	52	A23	N3-C4-N9	4.60	135.00	127.17
5	5	52	A23	C2-N3-C4	3.69	120.84	111.83
5	5	52	A23	O2C-PC-O1C	3.52	121.25	109.89
5	5	52	A23	C4-C5-N7	-3.47	106.62	110.58
5	5	52	A23	N3-C2-N1	-3.17	123.78	128.58
5	5	52	A23	C5'-C4'-C3'	2.88	123.96	114.38
5	5	52	A23	C4-N9-C8	2.62	108.49	105.74
5	5	52	A23	C5-N7-C8	2.53	107.43	103.45
5	5	52	A23	O3'-PC-O1C	-2.46	109.25	115.76
5	5	52	A23	O2'-PC-O1C	-2.46	109.26	115.76
5	5	52	A23	C2'-C1'-N9	-2.38	109.83	113.75
5	5	52	A23	C6-C5-N7	2.07	136.09	132.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	5	52	A23	3	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

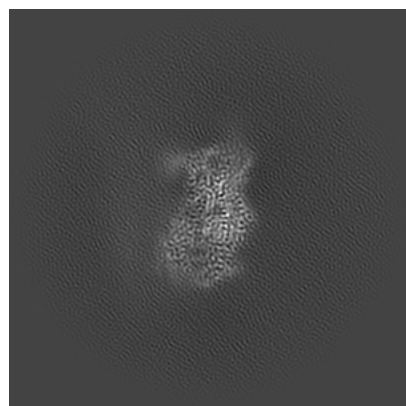
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34905. These allow visual inspection of the internal detail of the map and identification of artifacts.

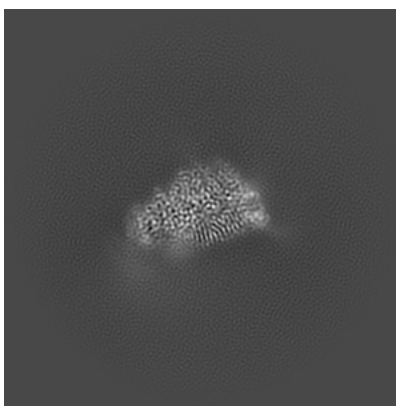
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

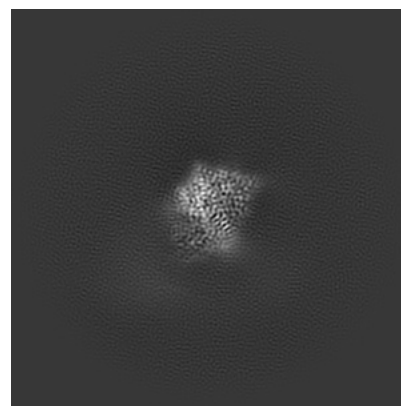
6.1.1 Primary map



X

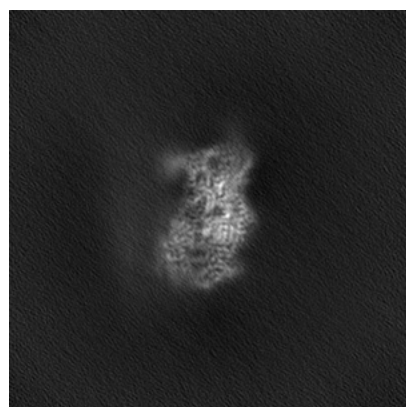


Y

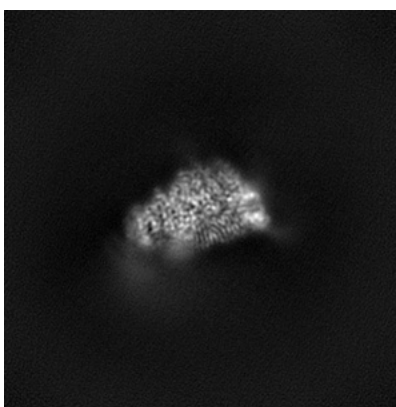


Z

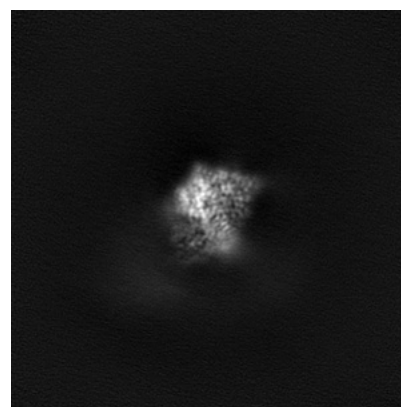
6.1.2 Raw map



X



Y

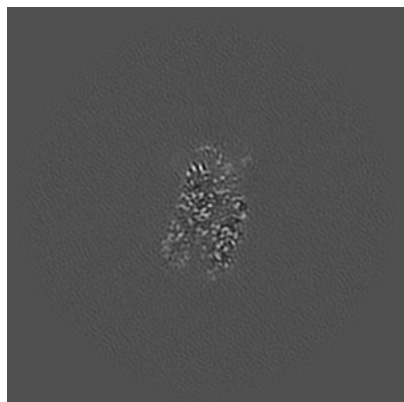


Z

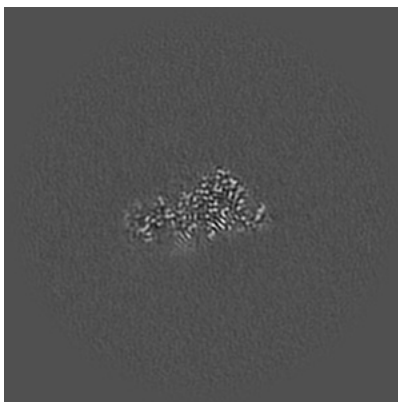
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

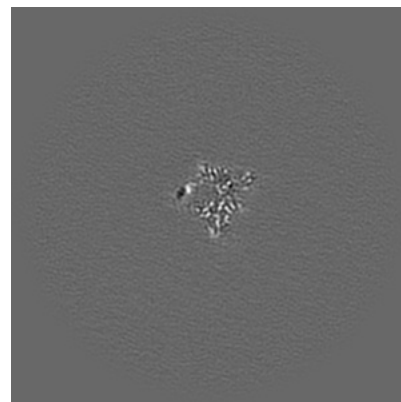
6.2.1 Primary map



X Index: 140

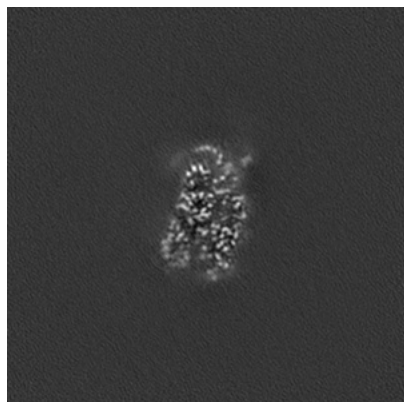


Y Index: 140

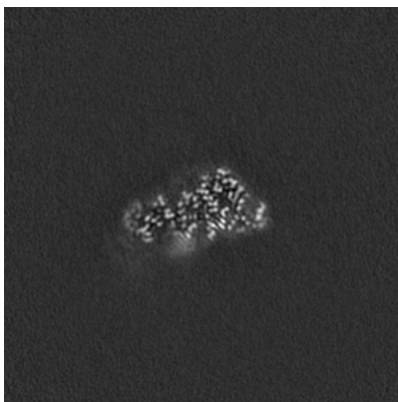


Z Index: 140

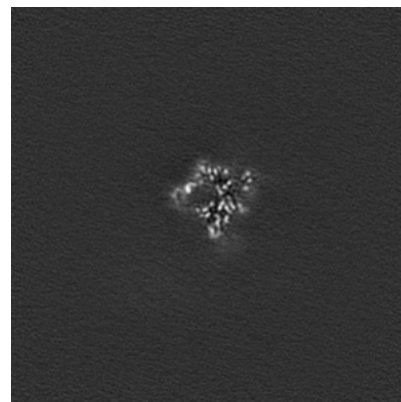
6.2.2 Raw map



X Index: 140



Y Index: 140

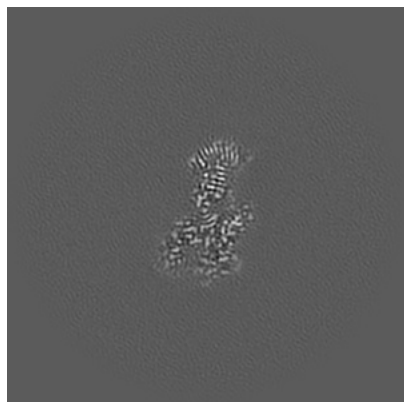


Z Index: 140

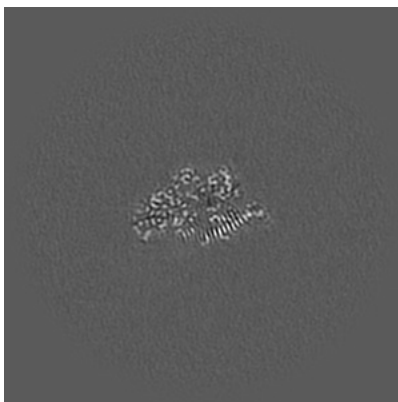
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

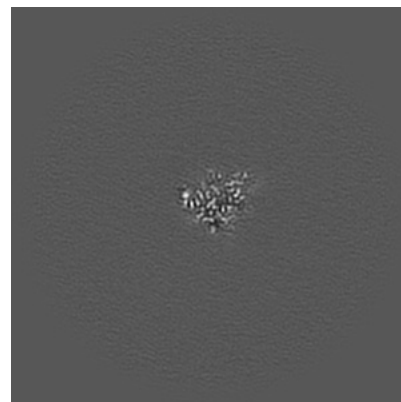
6.3.1 Primary map



X Index: 133

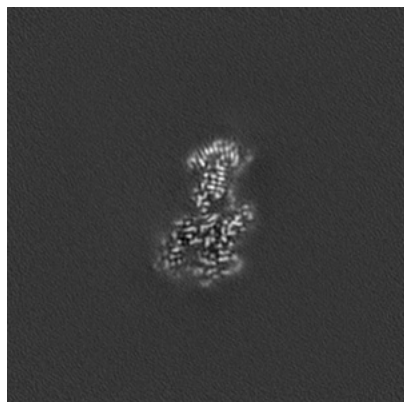


Y Index: 147

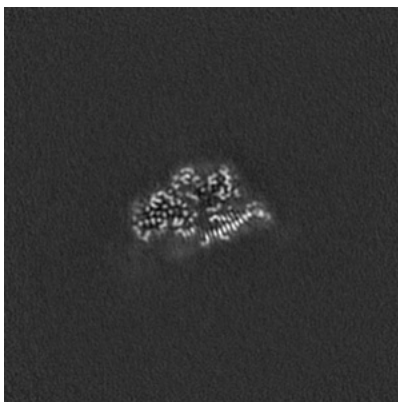


Z Index: 147

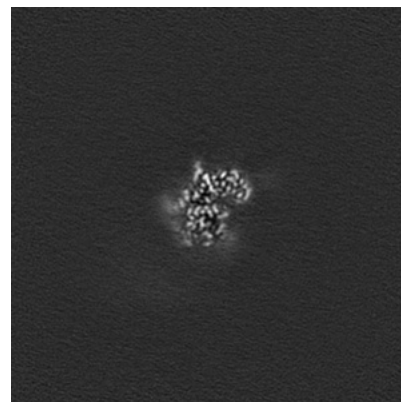
6.3.2 Raw map



X Index: 133



Y Index: 147

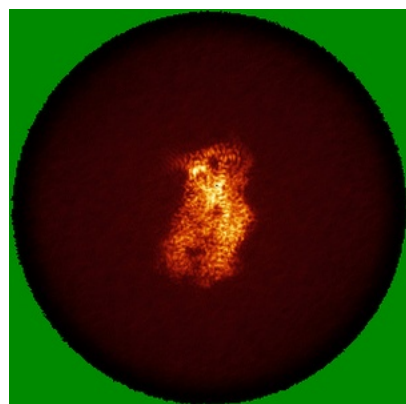


Z Index: 124

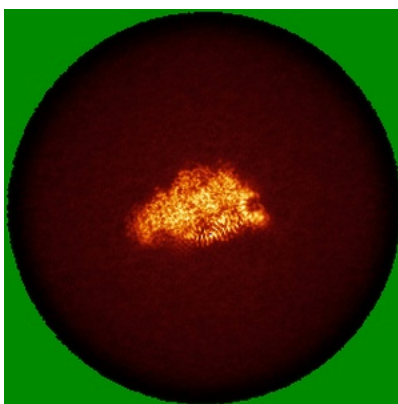
The images above show the largest variance slices of the map in three orthogonal directions.

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

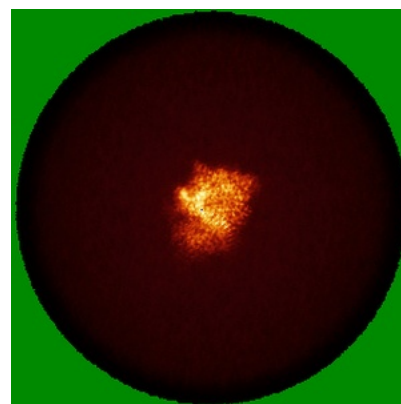
6.4.1 Primary map



X

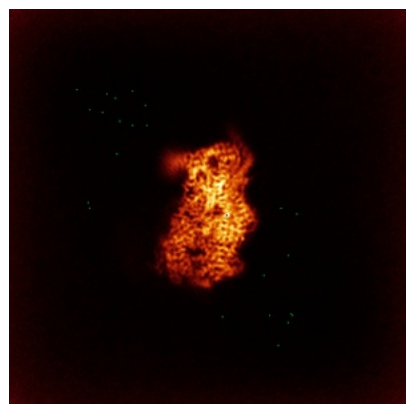


Y

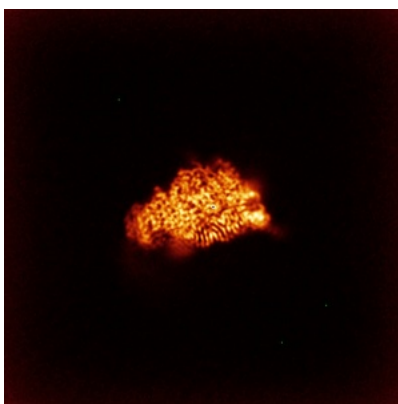


Z

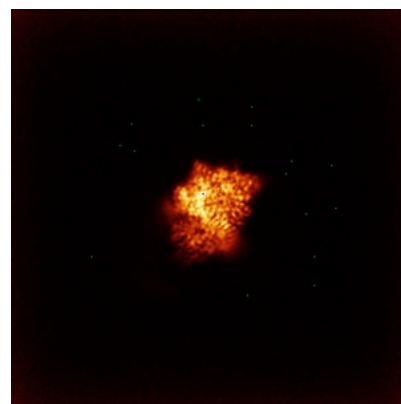
6.4.2 Raw map



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



X



Y



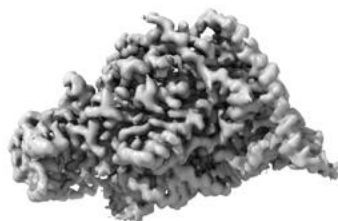
Z

The images above show the 3D surface view of the map at the recommended contour level 0.6. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

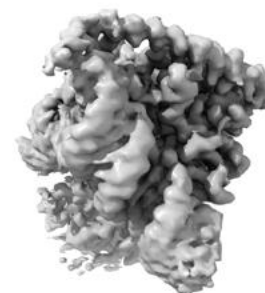
6.5.2 Raw map



X



Y



Z

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.

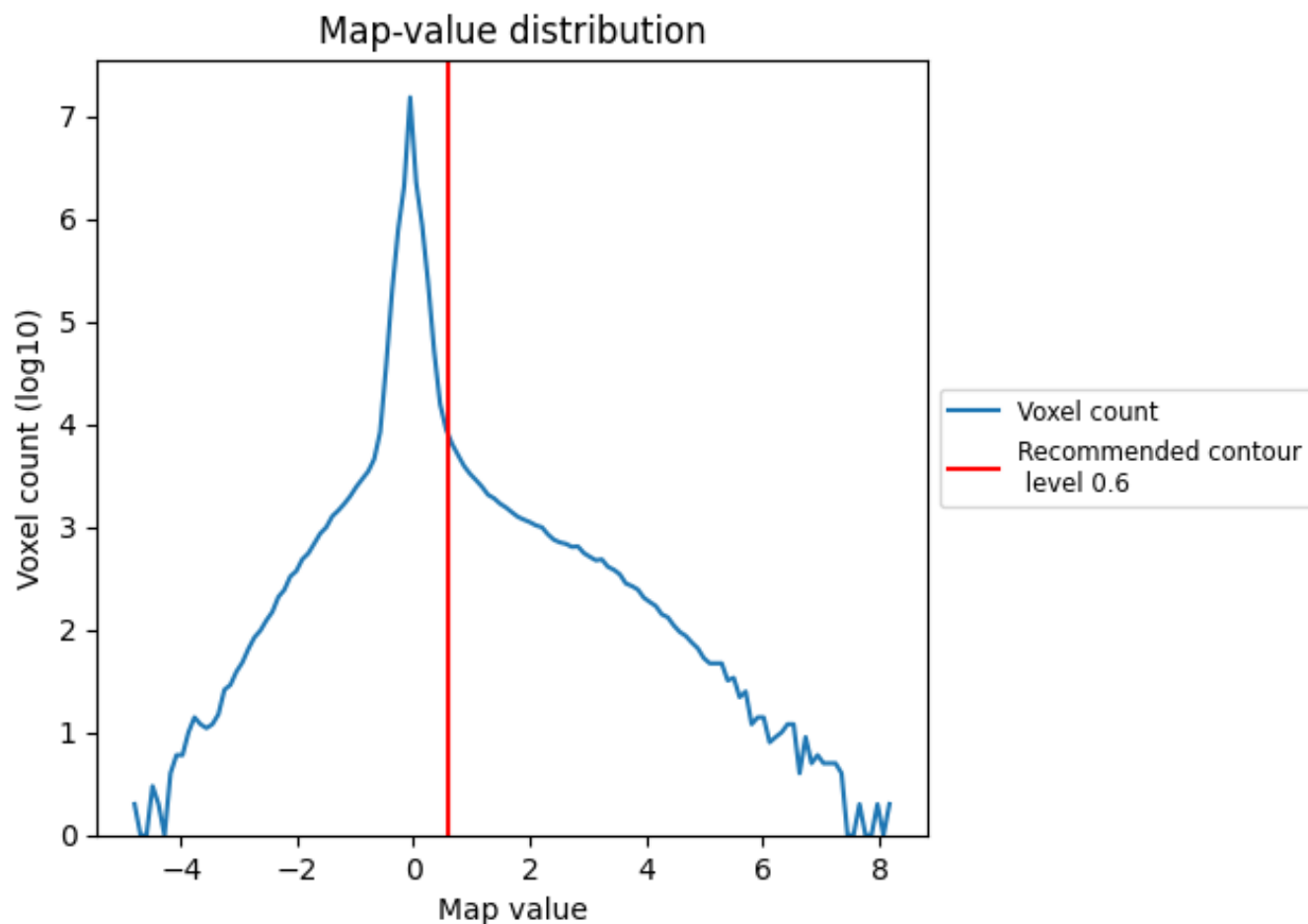
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

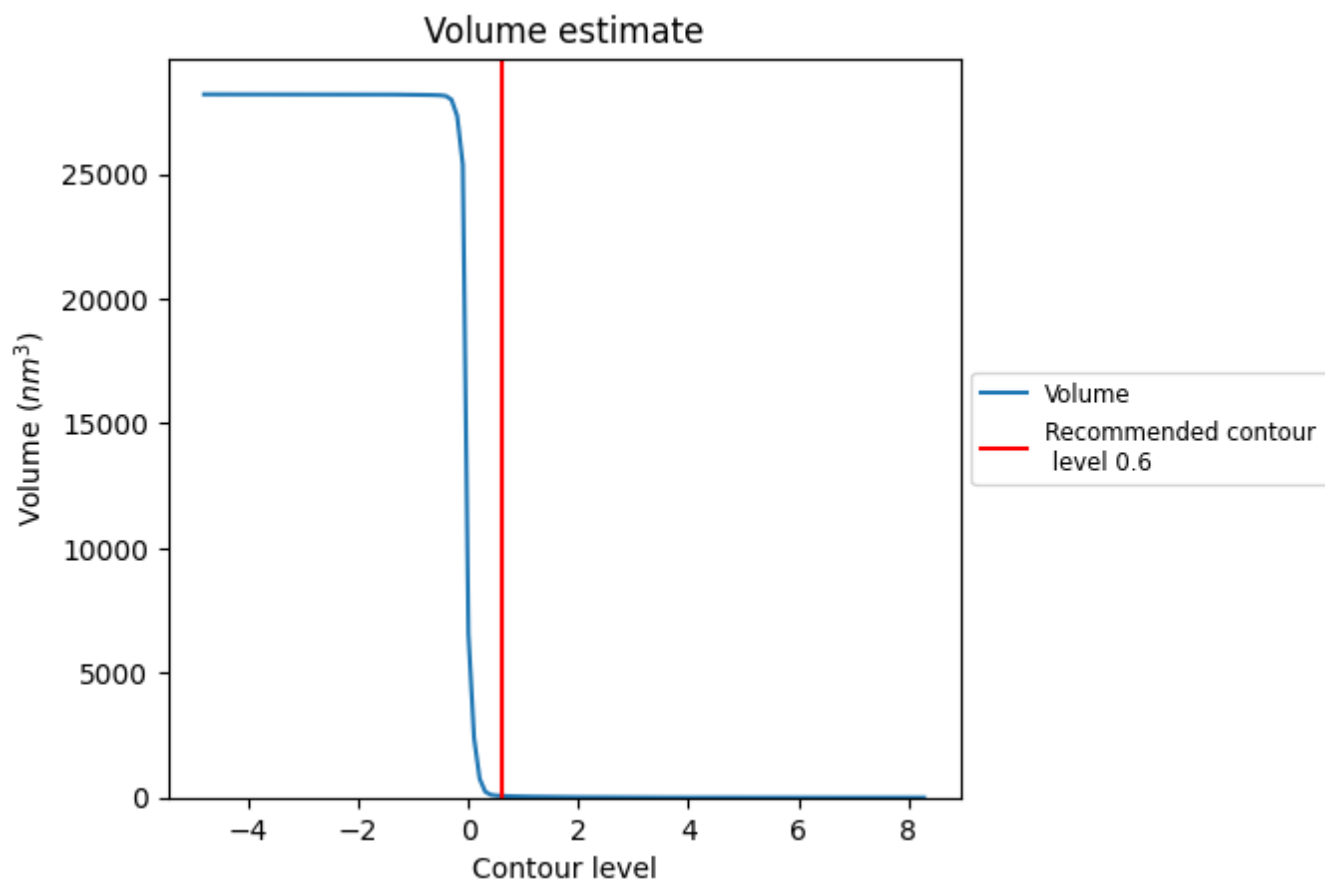
This section contains the results of statistical analysis of the map.

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.

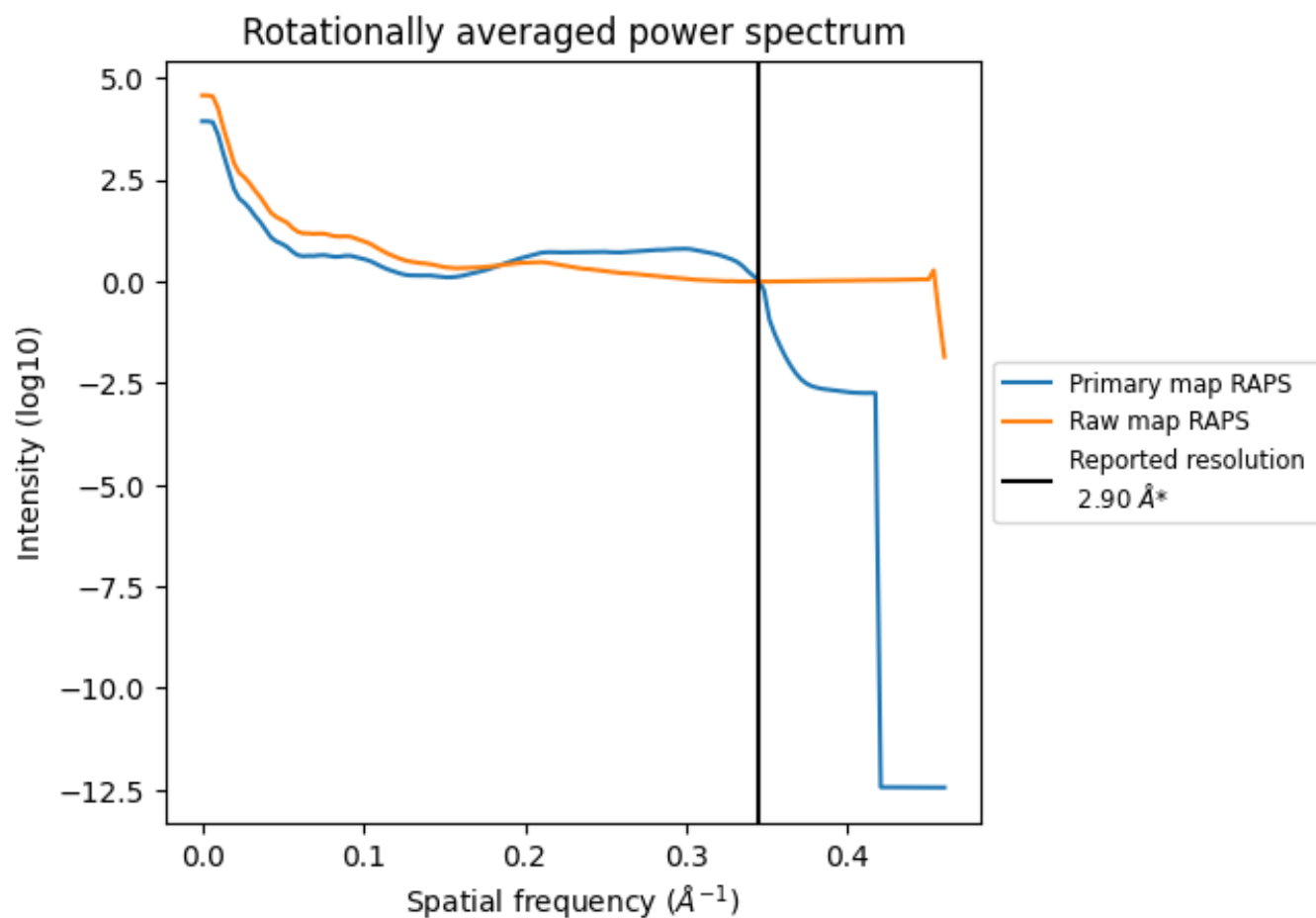
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 69 nm^3 ; this corresponds to an approximate mass of 62 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 ⓘ

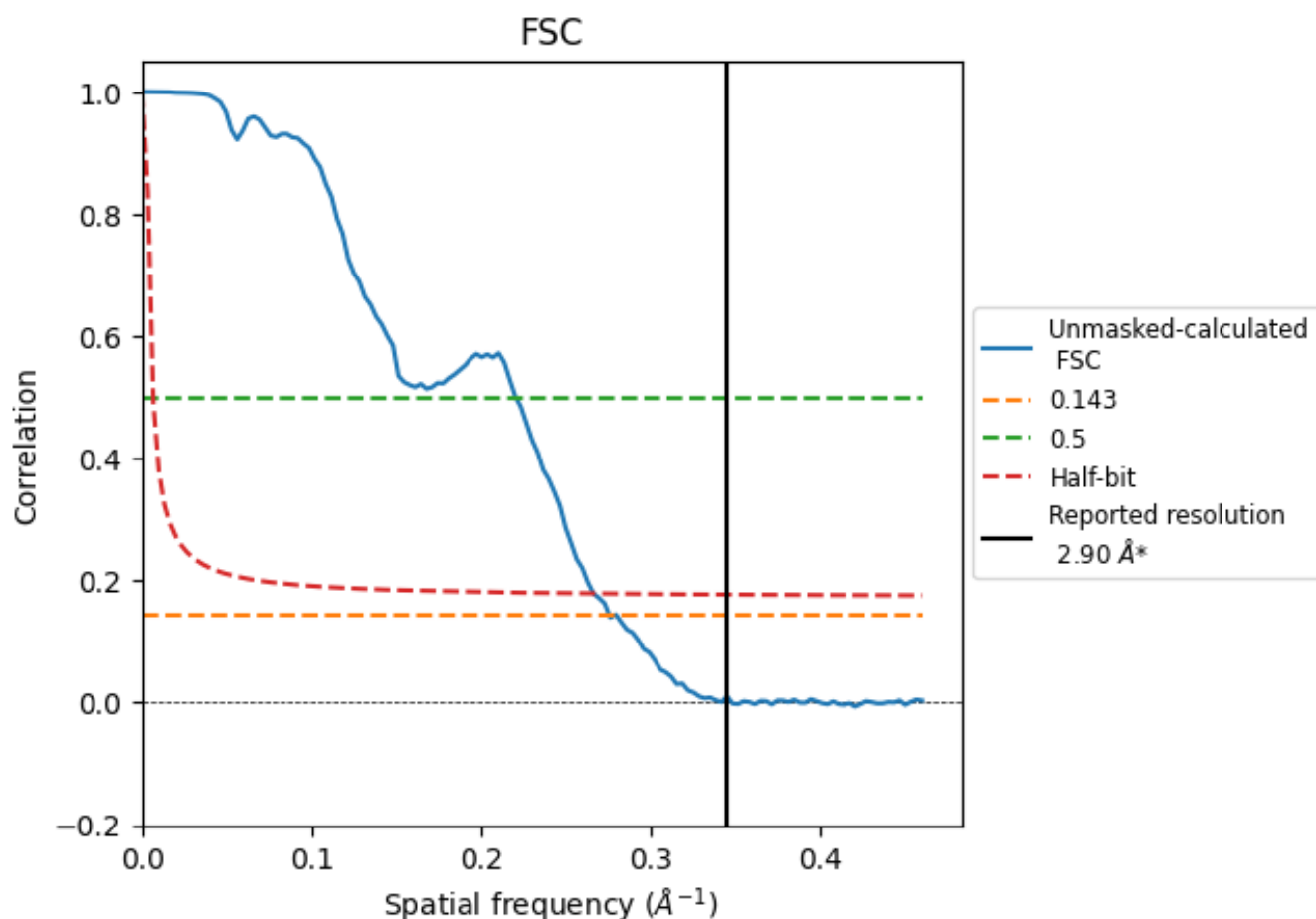


*Reported resolution corresponds to spatial frequency of 0.345 Å⁻¹

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 0.345 Å⁻¹

8.2 Resolution estimates [i](#)

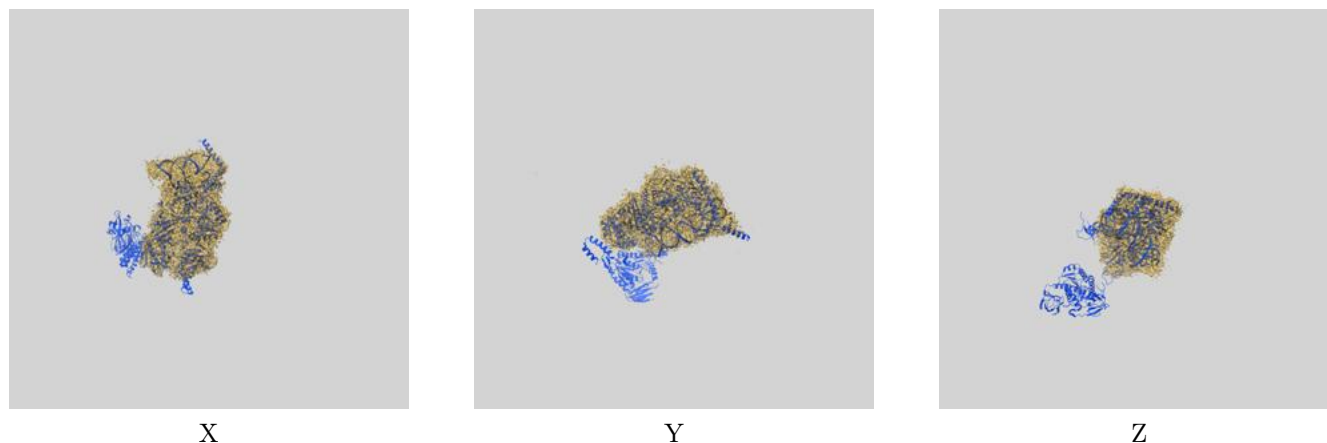
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.90	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.63	4.54	3.76

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.63 differs from the reported value 2.9 by more than 10 %

9 Map-model fit [i](#)

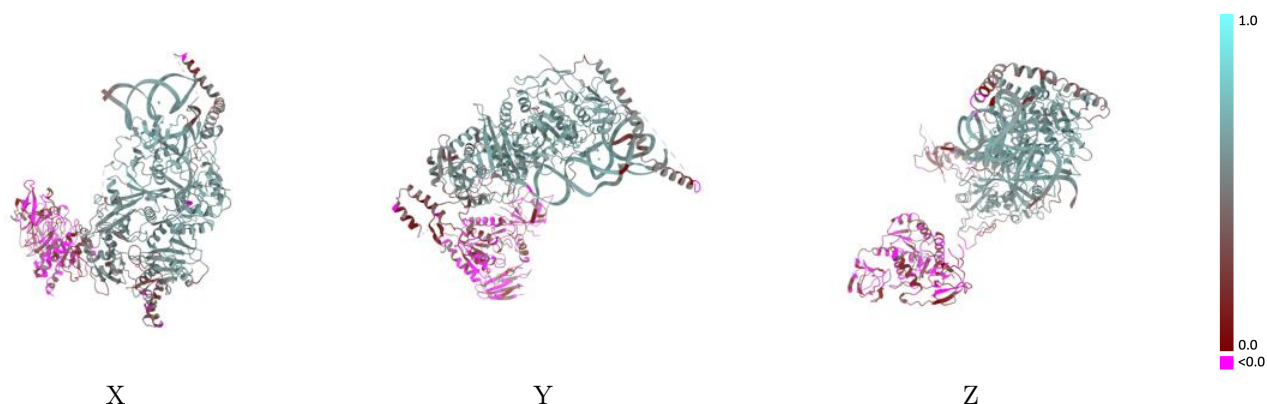
This section contains information regarding the fit between EMDB map EMD-34905 and PDB model 8HMZ. Per-residue inclusion information can be found in [section 3](#) on [page 9](#).

9.1 Map-model overlay [i](#)



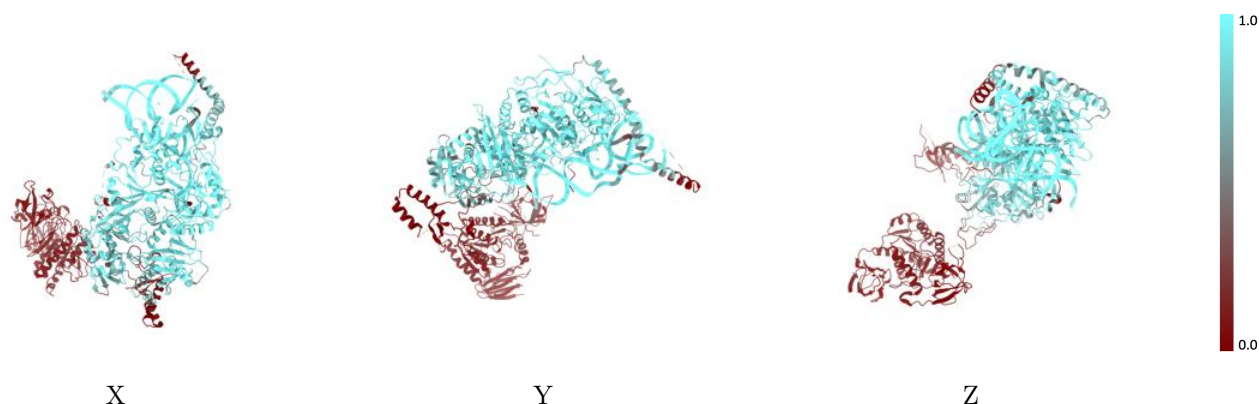
The images above show the 3D surface view of the map at the recommended contour level 0.6 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



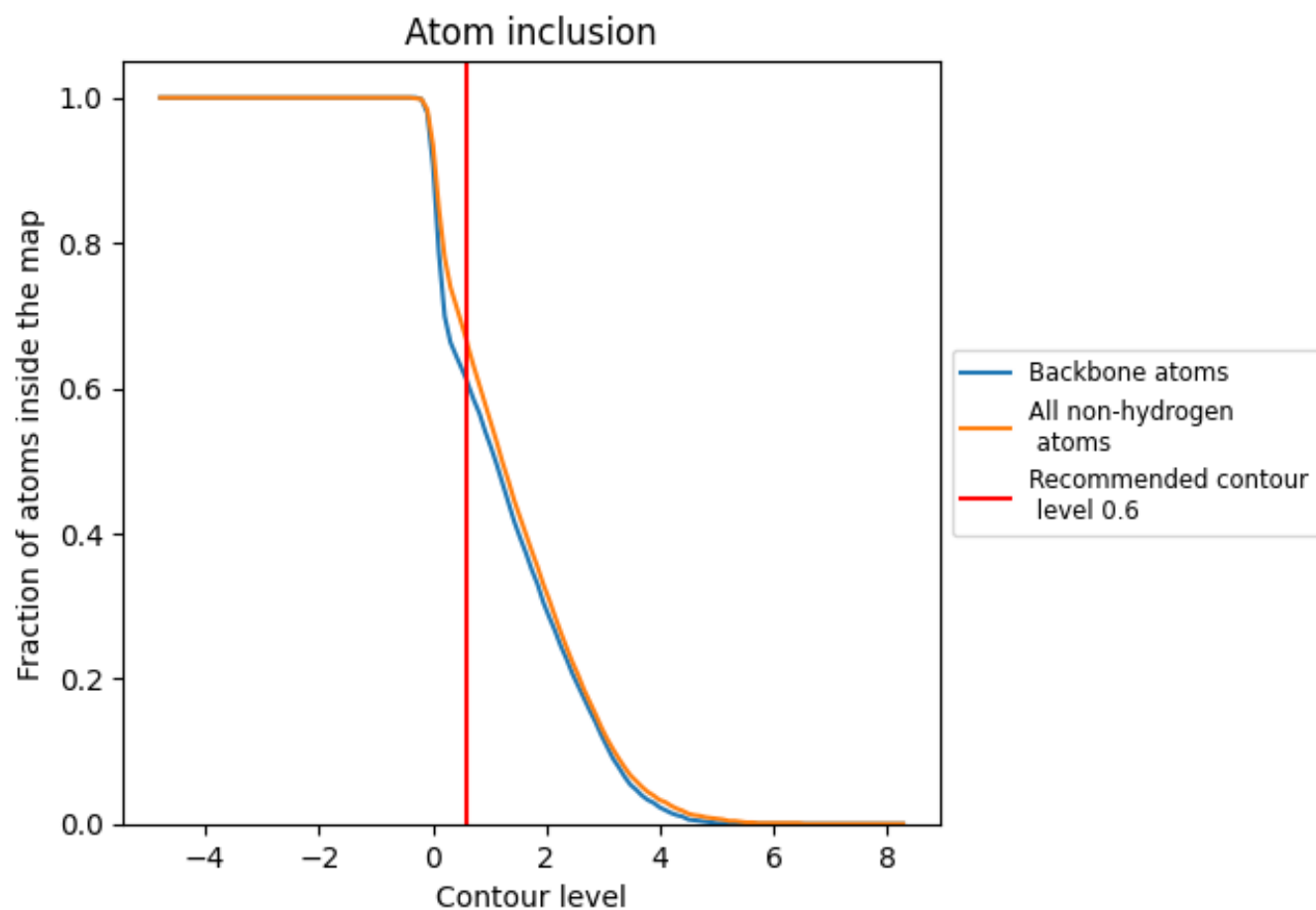
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.6).

9.4 Atom inclusion [i](#)



At the recommended contour level, 61% of all backbone atoms, 66% 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 (0.6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6650	0.4470
3	0.9710	0.5890
5	0.9010	0.5460
A	0.5410	0.4260
B	0.8720	0.5570
C	0.8960	0.5830
D	0.8280	0.5150
E	0.0000	0.0420

1.0

0.0

<0.0