Full wwPDB EM Validation Report (i)

Aug 31, 2026 – 05:43 PM EDT

PDB ID	:	12BJ / pdb_000012bj
EMDB ID	:	EMD-76280
Title	:	Moorella thermoacetica LarC (conformation 2)
Authors	:	Kim, K.H.; Wolfe, R.
Deposited on	:	2026-03-25
Resolution	:	3.85 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 (i) 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 (i)) were used in the production of this report:

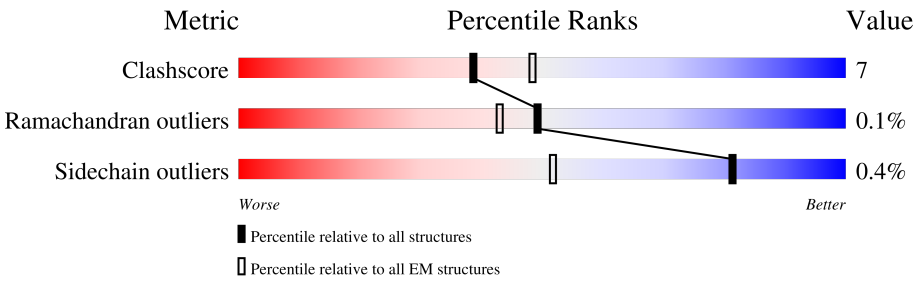
EMDB validation analysis	:	0.0.1.dev133
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

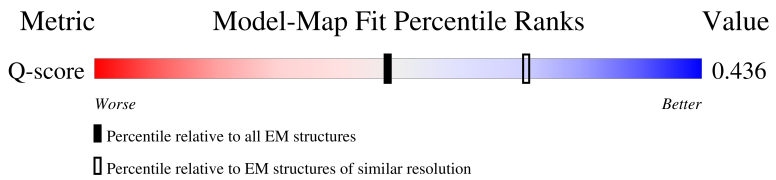
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.85 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	8989 (3.35 - 4.35)

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 <=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	414	14% 74% 18% 8%
1	B	414	7% 77% 18% •
1	C	414	• 83% 13% •
1	D	414	15% 79% 16% •

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Mol	Chain	Length	Quality of chain
1	E	414	5% 81% 15% •
1	F	414	17% 75% 17% 7%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 17592 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 Pyridinium-3,5-bisthiocarboxylic acid mononucleotide nickel insertion protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	E	400	2992	1908	515	555	14	0	0
1	A	382	2848	1813	487	535	13	0	0
1	B	396	2955	1882	508	551	14	0	0
1	C	400	2989	1904	515	556	14	0	0
1	F	383	2853	1816	488	536	13	0	0
1	D	396	2955	1882	508	551	14	0	0

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-17	MET	-	initiating methionine	UNP Q2RFJ5
E	-16	HIS	-	expression tag	UNP Q2RFJ5
E	-15	HIS	-	expression tag	UNP Q2RFJ5
E	-14	HIS	-	expression tag	UNP Q2RFJ5
E	-13	HIS	-	expression tag	UNP Q2RFJ5
E	-12	HIS	-	expression tag	UNP Q2RFJ5
E	-11	HIS	-	expression tag	UNP Q2RFJ5
E	-10	SER	-	expression tag	UNP Q2RFJ5
E	-9	GLY	-	expression tag	UNP Q2RFJ5
E	-8	THR	-	expression tag	UNP Q2RFJ5
E	-7	GLU	-	expression tag	UNP Q2RFJ5
E	-6	ASN	-	expression tag	UNP Q2RFJ5
E	-5	LEU	-	expression tag	UNP Q2RFJ5
E	-4	TYR	-	expression tag	UNP Q2RFJ5
E	-3	PHE	-	expression tag	UNP Q2RFJ5
E	-2	GLN	-	expression tag	UNP Q2RFJ5
E	-1	GLY	-	expression tag	UNP Q2RFJ5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	0	HIS	-	expression tag	UNP Q2RFJ5
A	-17	MET	-	initiating methionine	UNP Q2RFJ5
A	-16	HIS	-	expression tag	UNP Q2RFJ5
A	-15	HIS	-	expression tag	UNP Q2RFJ5
A	-14	HIS	-	expression tag	UNP Q2RFJ5
A	-13	HIS	-	expression tag	UNP Q2RFJ5
A	-12	HIS	-	expression tag	UNP Q2RFJ5
A	-11	HIS	-	expression tag	UNP Q2RFJ5
A	-10	SER	-	expression tag	UNP Q2RFJ5
A	-9	GLY	-	expression tag	UNP Q2RFJ5
A	-8	THR	-	expression tag	UNP Q2RFJ5
A	-7	GLU	-	expression tag	UNP Q2RFJ5
A	-6	ASN	-	expression tag	UNP Q2RFJ5
A	-5	LEU	-	expression tag	UNP Q2RFJ5
A	-4	TYR	-	expression tag	UNP Q2RFJ5
A	-3	PHE	-	expression tag	UNP Q2RFJ5
A	-2	GLN	-	expression tag	UNP Q2RFJ5
A	-1	GLY	-	expression tag	UNP Q2RFJ5
A	0	HIS	-	expression tag	UNP Q2RFJ5
B	-17	MET	-	initiating methionine	UNP Q2RFJ5
B	-16	HIS	-	expression tag	UNP Q2RFJ5
B	-15	HIS	-	expression tag	UNP Q2RFJ5
B	-14	HIS	-	expression tag	UNP Q2RFJ5
B	-13	HIS	-	expression tag	UNP Q2RFJ5
B	-12	HIS	-	expression tag	UNP Q2RFJ5
B	-11	HIS	-	expression tag	UNP Q2RFJ5
B	-10	SER	-	expression tag	UNP Q2RFJ5
B	-9	GLY	-	expression tag	UNP Q2RFJ5
B	-8	THR	-	expression tag	UNP Q2RFJ5
B	-7	GLU	-	expression tag	UNP Q2RFJ5
B	-6	ASN	-	expression tag	UNP Q2RFJ5
B	-5	LEU	-	expression tag	UNP Q2RFJ5
B	-4	TYR	-	expression tag	UNP Q2RFJ5
B	-3	PHE	-	expression tag	UNP Q2RFJ5
B	-2	GLN	-	expression tag	UNP Q2RFJ5
B	-1	GLY	-	expression tag	UNP Q2RFJ5
B	0	HIS	-	expression tag	UNP Q2RFJ5
C	-17	MET	-	initiating methionine	UNP Q2RFJ5
C	-16	HIS	-	expression tag	UNP Q2RFJ5
C	-15	HIS	-	expression tag	UNP Q2RFJ5
C	-14	HIS	-	expression tag	UNP Q2RFJ5
C	-13	HIS	-	expression tag	UNP Q2RFJ5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-12	HIS	-	expression tag	UNP Q2RFJ5
C	-11	HIS	-	expression tag	UNP Q2RFJ5
C	-10	SER	-	expression tag	UNP Q2RFJ5
C	-9	GLY	-	expression tag	UNP Q2RFJ5
C	-8	THR	-	expression tag	UNP Q2RFJ5
C	-7	GLU	-	expression tag	UNP Q2RFJ5
C	-6	ASN	-	expression tag	UNP Q2RFJ5
C	-5	LEU	-	expression tag	UNP Q2RFJ5
C	-4	TYR	-	expression tag	UNP Q2RFJ5
C	-3	PHE	-	expression tag	UNP Q2RFJ5
C	-2	GLN	-	expression tag	UNP Q2RFJ5
C	-1	GLY	-	expression tag	UNP Q2RFJ5
C	0	HIS	-	expression tag	UNP Q2RFJ5
F	-17	MET	-	initiating methionine	UNP Q2RFJ5
F	-16	HIS	-	expression tag	UNP Q2RFJ5
F	-15	HIS	-	expression tag	UNP Q2RFJ5
F	-14	HIS	-	expression tag	UNP Q2RFJ5
F	-13	HIS	-	expression tag	UNP Q2RFJ5
F	-12	HIS	-	expression tag	UNP Q2RFJ5
F	-11	HIS	-	expression tag	UNP Q2RFJ5
F	-10	SER	-	expression tag	UNP Q2RFJ5
F	-9	GLY	-	expression tag	UNP Q2RFJ5
F	-8	THR	-	expression tag	UNP Q2RFJ5
F	-7	GLU	-	expression tag	UNP Q2RFJ5
F	-6	ASN	-	expression tag	UNP Q2RFJ5
F	-5	LEU	-	expression tag	UNP Q2RFJ5
F	-4	TYR	-	expression tag	UNP Q2RFJ5
F	-3	PHE	-	expression tag	UNP Q2RFJ5
F	-2	GLN	-	expression tag	UNP Q2RFJ5
F	-1	GLY	-	expression tag	UNP Q2RFJ5
F	0	HIS	-	expression tag	UNP Q2RFJ5
D	-17	MET	-	initiating methionine	UNP Q2RFJ5
D	-16	HIS	-	expression tag	UNP Q2RFJ5
D	-15	HIS	-	expression tag	UNP Q2RFJ5
D	-14	HIS	-	expression tag	UNP Q2RFJ5
D	-13	HIS	-	expression tag	UNP Q2RFJ5
D	-12	HIS	-	expression tag	UNP Q2RFJ5
D	-11	HIS	-	expression tag	UNP Q2RFJ5
D	-10	SER	-	expression tag	UNP Q2RFJ5
D	-9	GLY	-	expression tag	UNP Q2RFJ5
D	-8	THR	-	expression tag	UNP Q2RFJ5
D	-7	GLU	-	expression tag	UNP Q2RFJ5

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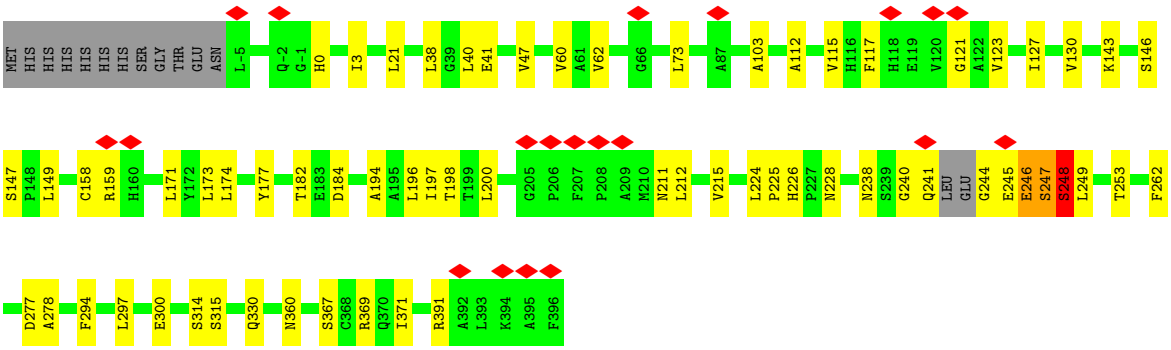
Chain	Residue	Modelled	Actual	Comment	Reference
D	-6	ASN	-	expression tag	UNP Q2RFJ5
D	-5	LEU	-	expression tag	UNP Q2RFJ5
D	-4	TYR	-	expression tag	UNP Q2RFJ5
D	-3	PHE	-	expression tag	UNP Q2RFJ5
D	-2	GLN	-	expression tag	UNP Q2RFJ5
D	-1	GLY	-	expression tag	UNP Q2RFJ5
D	0	HIS	-	expression tag	UNP Q2RFJ5

3 Residue-property plots (i)

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.

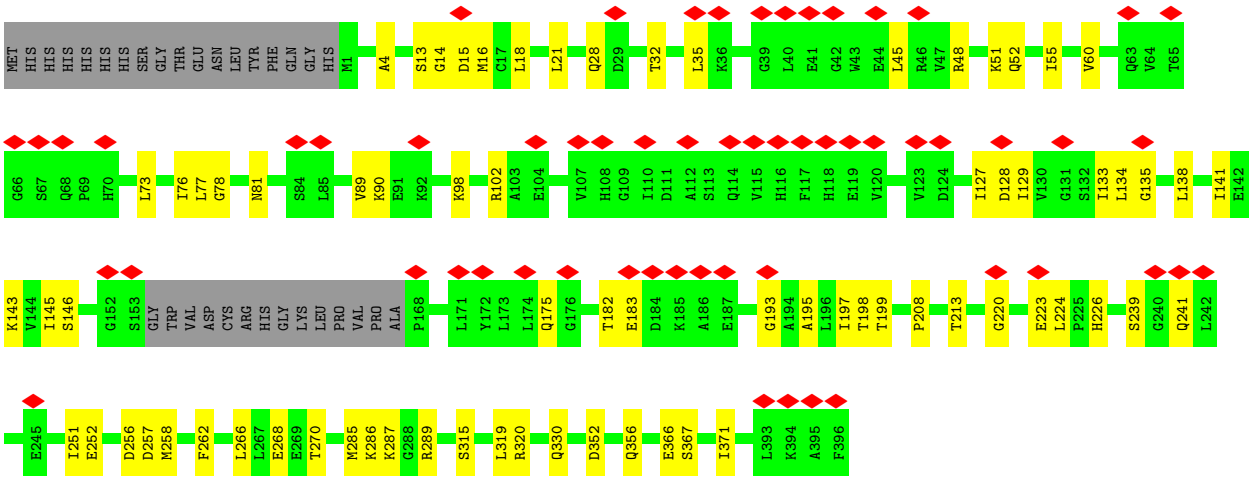
• Molecule 1: Pyridinium-3,5-bisthiocarboxylic acid mononucleotide nickel insertion protein

Chain E:



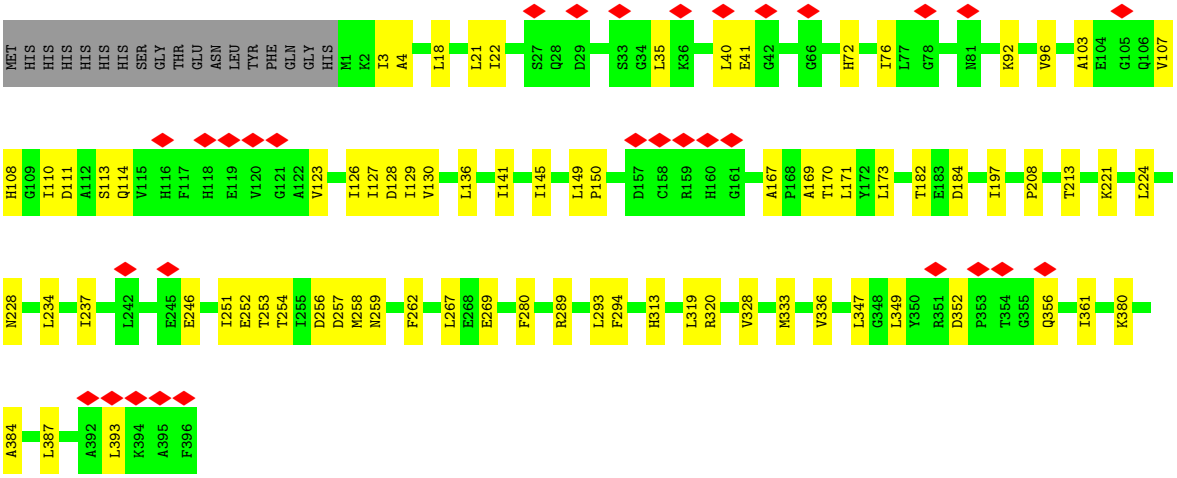
• Molecule 1: Pyridinium-3,5-bisthiocarboxylic acid mononucleotide nickel insertion protein

Chain A:



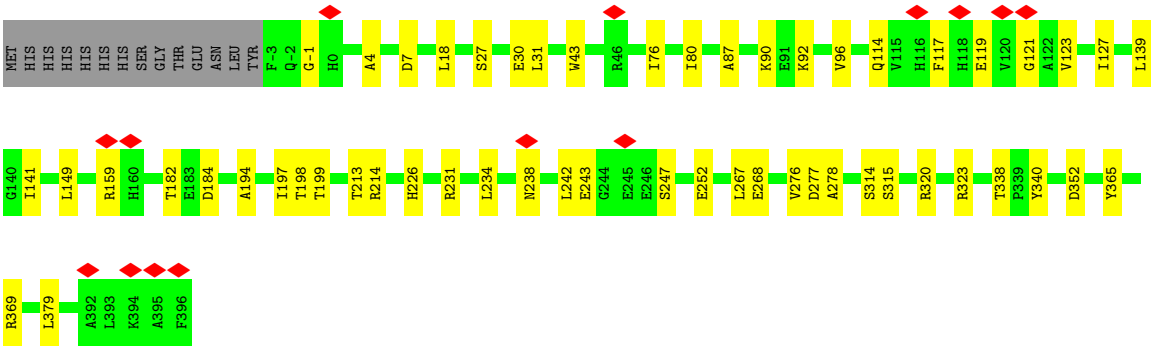
• Molecule 1: Pyridinium-3,5-bisthiocarboxylic acid mononucleotide nickel insertion protein

Chain B:



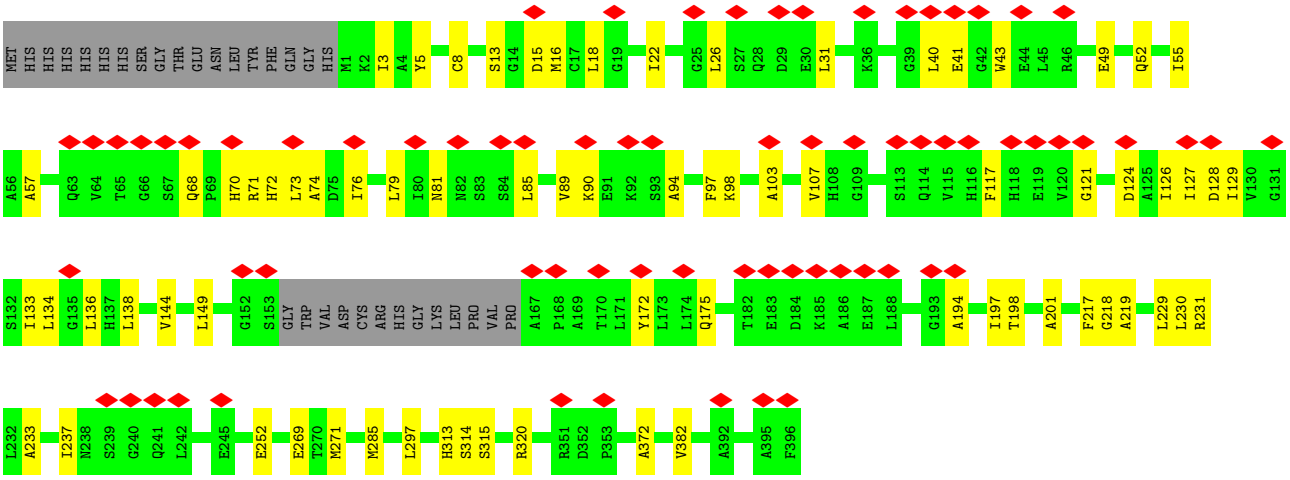
• Molecule 1: Pyridinium-3,5-bisthiocarboxylic acid mononucleotide nickel insertion protein

Chain C:



• Molecule 1: Pyridinium-3,5-bisthiocarboxylic acid mononucleotide nickel insertion protein

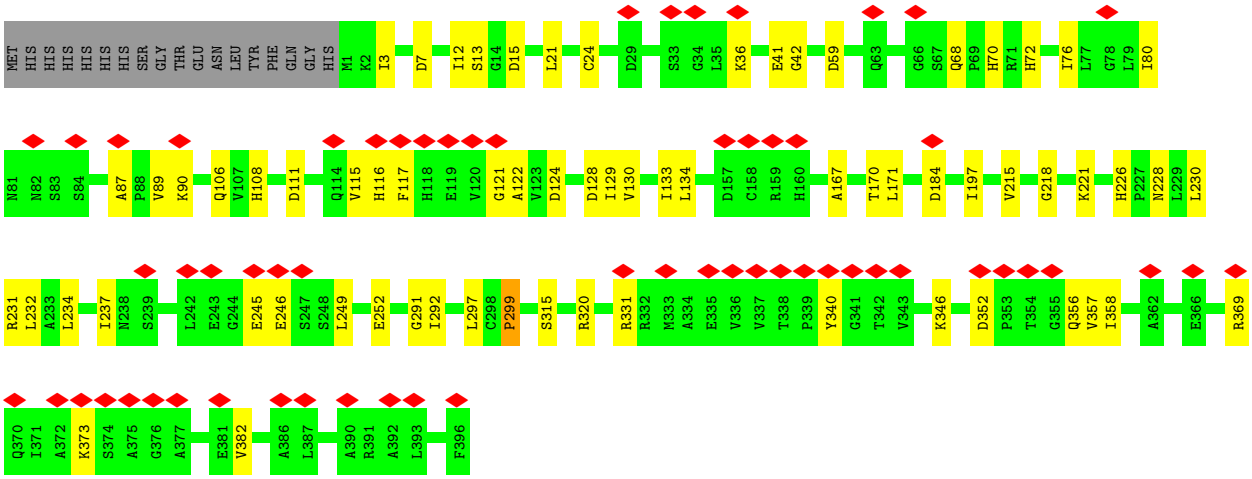
Chain F:



• Molecule 1: Pyridinium-3,5-bisthiocarboxylic acid mononucleotide nickel insertion protein

Chain D:

15%
79%
16%


MET HIS HIS HIS HIS HIS HIS SER GLY THR GLU ASN LEU TYR PHE GLN GLY HIS M1 K2 I3 D7 I12 S13 G14 D15 L21 C24 D29 S33 G34 L35 K36 E41 G42 D59 Q63 G66 S67 Q68 F69 H70 R71 H72 I76 L77 G78 L79 I80
N81 N82 S83 S84 A87 P88 V89 K90 Q106 V107 H108 D111 Q114 V115 H116 F117 H118 E119 V120 G121 A122 V123 D124 D128 I129 V130 I133 L134 D157 C158 R159 H160 A167 T170 L171 D184 I197 V215 G218 K221 H226 P227 N228 L229 L230
R231 L232 A233 L234 I237 N238 S239 L242 E243 G244 E245 E246 S247 S248 L249 E252 G291 I292 L297 C298 P299 S315 R320 R331 R332 M333 A334 E335 V336 V337 T338 P339 Y340 G341 T342 V343 K346 D352 P353 T354 G355 Q356 V357 I358 A362 E366 R369
Q370 I371 A372 K373 S374 A375 G376 A377 E381 V382 A386 L387 A390 R391 A392 L393 F396

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	142648	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 ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.034	Depositor
Minimum map value	-0.670	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.12	Depositor
Map size (Å)	325.9008, 325.9008, 325.9008	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8487, 0.8487, 0.8487	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	0.13	0/2902	0.36	0/3939
1	B	0.12	0/3015	0.31	0/4097
1	C	0.12	0/3051	0.29	0/4145
1	D	0.19	1/3015 (0.0%)	0.44	4/4097 (0.1%)
1	E	0.20	0/3054	0.39	4/4148 (0.1%)
1	F	0.12	0/2907	0.32	0/3947
All	All	0.15	1/17944 (0.0%)	0.36	8/24373 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	299	PRO	CG-CD	-8.14	1.23	1.50

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	299	PRO	N-CD-CG	-13.65	82.72	103.20
1	D	299	PRO	CA-CB-CG	-10.55	84.44	104.50
1	E	225	PRO	CA-N-CD	-8.55	100.03	112.00
1	E	225	PRO	N-CD-CG	-6.64	93.24	103.20
1	D	299	PRO	N-CA-CB	-6.38	95.73	103.45
1	D	299	PRO	CA-N-CD	-5.74	103.97	112.00
1	E	225	PRO	CB-CG-CD	-5.06	89.90	106.10
1	E	225	PRO	CA-CB-CG	-5.01	94.99	104.50

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	2848	0	2896	49	0
1	B	2955	0	3002	48	0
1	C	2989	0	3029	33	0
1	D	2955	0	3002	44	0
1	E	2992	0	3031	44	0
1	F	2853	0	2900	49	0
All	All	17592	0	17860	257	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (257) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:MET:HE1	1:A:220:GLY:HA2	1.58	0.84
1:E:0:HIS:H	1:E:238:ASN:HA	1.47	0.79
1:C:96:VAL:HG22	1:C:199:THR:HG21	1.66	0.75
1:C:4:ALA:HB2	1:C:141:ILE:HD12	1.69	0.74
1:E:158:CYS:SG	1:E:159:ARG:N	2.62	0.72
1:C:213:THR:HG22	1:C:214:ARG:HG2	1.72	0.72
1:C:114:GLN:N	1:C:114:GLN:OE1	2.24	0.71
1:D:3:ILE:HD13	1:D:237:ILE:HD11	1.74	0.70
1:F:13:SER:OG	1:F:15:ASP:OD1	2.11	0.67
1:A:35:LEU:HD21	1:A:127:ILE:HD11	1.76	0.66
1:C:123:VAL:O	1:C:127:ILE:HG22	1.95	0.66
1:F:89:VAL:HG11	1:F:134:LEU:HD22	1.76	0.66
1:A:258:MET:HE2	1:A:262:PHE:HB2	1.79	0.65
1:E:123:VAL:O	1:E:127:ILE:HG22	1.97	0.64
1:A:73:LEU:O	1:A:77:LEU:HD12	1.98	0.64
1:F:3:ILE:HD13	1:F:237:ILE:HD11	1.80	0.64
1:B:4:ALA:HB2	1:B:141:ILE:HD12	1.81	0.63
1:E:367:SER:O	1:E:371:ILE:HG12	1.99	0.63
1:A:175:GLN:OE1	1:A:175:GLN:N	2.25	0.63
1:A:252:GLU:OE2	1:A:320:ARG:NH2	2.34	0.61
1:D:245:GLU:O	1:D:246:GLU:HG3	2.00	0.61
1:C:182:THR:HG22	1:C:184:ASP:H	1.64	0.61
1:A:224:LEU:HD22	1:A:226:HIS:HE1	1.65	0.61
1:A:256:ASP:OD2	1:A:289:ARG:NH2	2.33	0.61
1:A:128:ASP:OD1	1:A:129:ILE:N	2.34	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:-1:GLY:HA2	1:C:238:ASN:HA	1.83	0.61
1:E:38:LEU:HD11	1:E:130:VAL:HG11	1.83	0.60
1:B:336:VAL:HG22	1:B:393:LEU:HD23	1.82	0.60
1:F:49:GLU:OE2	1:F:49:GLU:N	2.34	0.60
1:E:315:SER:O	1:E:315:SER:OG	2.18	0.60
1:D:129:ILE:O	1:D:133:ILE:HG12	2.01	0.60
1:B:128:ASP:OD1	1:B:129:ILE:N	2.35	0.60
1:F:194:ALA:O	1:F:198:THR:HG22	2.01	0.60
1:D:7:ASP:HB3	1:D:231:ARG:HB2	1.83	0.60
1:F:81:ASN:O	1:F:90:LYS:NZ	2.30	0.59
1:F:314:SER:OG	1:F:315:SER:N	2.35	0.59
1:D:130:VAL:O	1:D:134:LEU:HG	2.02	0.59
1:E:182:THR:HG22	1:E:184:ASP:H	1.68	0.59
1:F:117:PHE:HB3	1:F:121:GLY:HA3	1.84	0.58
1:A:73:LEU:O	1:A:76:ILE:HG12	2.04	0.58
1:A:21:LEU:HD13	1:A:135:GLY:HA3	1.85	0.58
1:C:314:SER:OG	1:C:315:SER:N	2.35	0.58
1:E:103:ALA:HB2	1:E:173:LEU:HD13	1.84	0.58
1:E:369:ARG:NH2	1:C:268:GLU:OE1	2.37	0.58
1:B:256:ASP:OD2	1:B:289:ARG:NH2	2.37	0.58
1:F:252:GLU:OE2	1:F:320:ARG:NH2	2.37	0.57
1:B:126:ILE:O	1:B:130:VAL:HG23	2.05	0.57
1:F:18:LEU:HD11	1:F:128:ASP:HA	1.85	0.57
1:F:68:GLN:OE1	1:F:71:ARG:NH2	2.38	0.57
1:D:13:SER:OG	1:D:15:ASP:OD1	2.20	0.57
1:A:195:ALA:O	1:A:199:THR:HG22	2.05	0.56
1:C:139:LEU:HD13	1:C:234:LEU:HD11	1.86	0.56
1:C:338:THR:HG23	1:C:340:TYR:H	1.69	0.56
1:E:146:SER:OG	1:E:147:SER:N	2.37	0.56
1:E:0:HIS:NE2	1:E:330:GLN:OE1	2.38	0.56
1:A:51:LYS:NZ	1:B:108:HIS:O	2.37	0.56
1:B:40:LEU:HD21	1:B:123:VAL:HB	1.88	0.56
1:A:193:GLY:O	1:A:197:ILE:HG22	2.05	0.56
1:E:247:SER:HB3	1:E:300:GLU:OE1	2.05	0.56
1:A:77:LEU:O	1:A:81:ASN:ND2	2.39	0.56
1:D:106:GLN:O	1:D:106:GLN:NE2	2.39	0.56
1:A:18:LEU:HB2	1:A:60:VAL:HG11	1.88	0.55
1:E:246:GLU:O	1:E:247:SER:C	2.48	0.55
1:B:111:ASP:HB3	1:B:114:GLN:HE21	1.72	0.55
1:E:117:PHE:HB3	1:E:121:GLY:HA3	1.88	0.55
1:F:134:LEU:O	1:F:138:LEU:HG	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:97:PHE:HE2	1:F:126:ILE:HD13	1.70	0.55
1:B:40:LEU:HD12	1:B:41:GLU:H	1.72	0.55
1:C:92:LYS:O	1:C:96:VAL:HG23	2.07	0.55
1:A:4:ALA:HB2	1:A:141:ILE:HD12	1.88	0.55
1:B:224:LEU:HD12	1:B:228:ASN:HB3	1.90	0.54
1:B:333:MET:HE3	1:B:333:MET:HA	1.88	0.54
1:C:149:LEU:HD21	1:C:197:ILE:HG21	1.89	0.54
1:D:340:TYR:HD2	1:D:382:VAL:HG23	1.72	0.54
1:C:194:ALA:O	1:C:198:THR:HG23	2.08	0.54
1:D:15:ASP:OD1	1:D:15:ASP:N	2.37	0.53
1:D:68:GLN:HE22	1:D:122:ALA:HB1	1.73	0.53
1:A:14:GLY:O	1:A:18:LEU:HG	2.08	0.53
1:B:18:LEU:O	1:B:22:ILE:HG23	2.08	0.53
1:F:269:GLU:OE1	1:F:313:HIS:NE2	2.42	0.53
1:E:240:GLY:O	1:E:241:GLN:C	2.51	0.53
1:B:111:ASP:OD2	1:B:113:SER:OG	2.23	0.53
1:B:246:GLU:N	1:B:246:GLU:OE1	2.41	0.53
1:F:5:TYR:HB3	1:F:233:ALA:HB3	1.91	0.53
1:E:184:ASP:OD2	1:E:226:HIS:NE2	2.40	0.53
1:D:331:ARG:HB3	1:D:346:LYS:HE3	1.91	0.53
1:A:98:LYS:O	1:A:98:LYS:NZ	2.37	0.52
1:F:41:GLU:O	1:F:68:GLN:NE2	2.42	0.52
1:D:72:HIS:O	1:D:76:ILE:HG12	2.10	0.52
1:C:141:ILE:HD11	1:C:234:LEU:HD22	1.92	0.51
1:B:258:MET:HE2	1:B:262:PHE:HB2	1.91	0.51
1:C:365:TYR:OH	1:C:369:ARG:NH2	2.43	0.51
1:F:18:LEU:O	1:F:22:ILE:HD12	2.11	0.51
1:F:70:HIS:C	1:F:70:HIS:CD2	2.89	0.51
1:D:12:ILE:HD13	1:D:197:ILE:HB	1.91	0.51
1:A:15:ASP:N	1:A:15:ASP:OD1	2.43	0.51
1:D:184:ASP:OD1	1:D:226:HIS:NE2	2.39	0.51
1:A:315:SER:O	1:A:315:SER:OG	2.22	0.51
1:F:89:VAL:HG13	1:F:133:ILE:HG22	1.92	0.51
1:F:73:LEU:HD12	1:F:74:ALA:N	2.26	0.51
1:A:266:LEU:O	1:A:270:THR:HG23	2.11	0.51
1:E:314:SER:OG	1:E:315:SER:N	2.44	0.50
1:B:149:LEU:HD21	1:B:197:ILE:HG21	1.94	0.50
1:B:251:ILE:HG23	1:B:319:LEU:HD13	1.93	0.50
1:B:352:ASP:OD1	1:B:356:GLN:N	2.45	0.50
1:F:94:ALA:O	1:F:98:LYS:HG3	2.11	0.50
1:F:97:PHE:CE2	1:F:126:ILE:HD13	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:360:ASN:OD1	1:B:380:LYS:NZ	2.45	0.50
1:C:184:ASP:OD2	1:C:226:HIS:NE2	2.38	0.50
1:D:76:ILE:O	1:D:80:ILE:HG12	2.11	0.50
1:A:146:SER:HB3	1:A:197:ILE:HD11	1.94	0.50
1:D:352:ASP:OD1	1:D:356:GLN:N	2.45	0.49
1:E:3:ILE:HG22	1:E:143:LYS:HB3	1.93	0.49
1:A:330:GLN:OE1	1:A:330:GLN:N	2.45	0.49
1:C:252:GLU:OE2	1:C:320:ARG:NH2	2.45	0.49
1:F:26:LEU:HD21	1:F:31:LEU:HD22	1.94	0.49
1:D:252:GLU:OE2	1:D:320:ARG:NH2	2.36	0.49
1:F:128:ASP:OD1	1:F:129:ILE:N	2.45	0.49
1:E:149:LEU:HD21	1:E:197:ILE:HG21	1.95	0.49
1:A:13:SER:OG	1:A:15:ASP:OD1	2.23	0.49
1:B:254:THR:HG22	1:B:293:LEU:HD12	1.94	0.49
1:D:117:PHE:HB3	1:D:121:GLY:HA3	1.94	0.49
1:A:352:ASP:OD1	1:A:356:GLN:N	2.45	0.48
1:C:87:ALA:HA	1:C:90:LYS:HD3	1.94	0.48
1:B:182:THR:HG23	1:B:184:ASP:H	1.78	0.48
1:E:171:LEU:HG	1:D:231:ARG:HH21	1.79	0.48
1:B:252:GLU:OE1	1:B:320:ARG:NH2	2.46	0.48
1:F:72:HIS:O	1:F:76:ILE:HG13	2.13	0.48
1:E:47:VAL:HG12	1:E:60:VAL:HG22	1.94	0.47
1:E:277:ASP:OD1	1:E:278:ALA:N	2.47	0.47
1:D:128:ASP:OD1	1:D:129:ILE:N	2.47	0.47
1:D:21:LEU:HD22	1:D:232:LEU:HD21	1.97	0.47
1:F:85:LEU:O	1:F:90:LYS:HE3	2.15	0.47
1:F:129:ILE:O	1:F:133:ILE:HG12	2.15	0.47
1:A:52:GLN:O	1:A:55:ILE:HG12	2.15	0.47
1:D:245:GLU:OE2	1:D:249:LEU:HD22	2.13	0.47
1:E:253:THR:HG23	1:E:294:PHE:HB3	1.96	0.47
1:E:391:ARG:HH22	1:B:384:ALA:HB1	1.78	0.47
1:A:81:ASN:HA	1:A:90:LYS:HD2	1.97	0.47
1:B:21:LEU:HD21	1:B:136:LEU:HD22	1.96	0.47
1:C:352:ASP:OD1	1:C:352:ASP:N	2.45	0.47
1:F:15:ASP:OD1	1:F:15:ASP:N	2.47	0.47
1:F:149:LEU:HD21	1:F:197:ILE:HG21	1.97	0.47
1:F:52:GLN:O	1:F:55:ILE:HG12	2.15	0.47
1:A:28:GLN:O	1:A:32:THR:HG23	2.16	0.46
1:A:367:SER:O	1:A:371:ILE:HG12	2.15	0.46
1:E:177:TYR:CD2	1:E:198:THR:HG22	2.50	0.46
1:F:103:ALA:O	1:F:107:VAL:HG22	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:THR:OG1	1:A:183:GLU:N	2.49	0.46
1:F:70:HIS:HD2	1:F:71:ARG:N	2.13	0.46
1:F:372:ALA:HB2	1:F:382:VAL:HG11	1.97	0.46
1:B:92:LYS:O	1:B:96:VAL:HG12	2.16	0.46
1:D:111:ASP:O	1:D:115:VAL:HG23	2.15	0.46
1:A:32:THR:HG22	1:A:45:LEU:HD12	1.98	0.46
1:B:253:THR:HG23	1:B:294:PHE:HB3	1.97	0.45
1:F:149:LEU:HD21	1:F:197:ILE:HD13	1.97	0.45
1:D:315:SER:O	1:D:315:SER:OG	2.33	0.45
1:A:257:ASP:N	1:A:257:ASP:OD1	2.48	0.45
1:B:167:ALA:O	1:B:170:THR:OG1	2.27	0.45
1:A:268:GLU:OE1	1:D:369:ARG:NH2	2.49	0.45
1:D:87:ALA:O	1:D:90:LYS:HG2	2.16	0.45
1:A:366:GLU:N	1:A:366:GLU:OE2	2.50	0.45
1:D:291:GLY:C	1:D:292:ILE:HD12	2.41	0.45
1:A:285:MET:HG3	1:A:286:LYS:H	1.81	0.45
1:F:40:LEU:HB3	1:F:43:TRP:HE1	1.82	0.45
1:B:103:ALA:O	1:B:107:VAL:HG23	2.16	0.45
1:F:57:ALA:HB1	1:F:219:ALA:HB2	1.99	0.45
1:C:7:ASP:HB3	1:C:231:ARG:HB3	1.99	0.44
1:C:276:VAL:HG12	1:C:277:ASP:OD1	2.18	0.44
1:D:89:VAL:HG22	1:D:133:ILE:HG22	1.99	0.44
1:D:357:VAL:C	1:D:358:ILE:HD13	2.42	0.44
1:E:40:LEU:HD12	1:E:41:GLU:H	1.82	0.44
1:E:262:PHE:CD1	1:C:379:LEU:HD23	2.53	0.44
1:B:169:ALA:O	1:B:173:LEU:HD13	2.17	0.44
1:D:24:CYS:SG	1:D:234:LEU:HD11	2.58	0.44
1:C:247:SER:OG	1:C:323:ARG:NE	2.47	0.44
1:F:136:LEU:HD23	1:F:136:LEU:HA	1.87	0.44
1:D:59:ASP:OD2	1:D:221:LYS:NZ	2.50	0.44
1:A:195:ALA:HA	1:A:198:THR:HG22	1.99	0.44
1:B:269:GLU:HG3	1:B:313:HIS:CE1	2.53	0.44
1:E:177:TYR:HD2	1:E:198:THR:HG22	1.83	0.44
1:A:145:ILE:HD12	1:A:208:PRO:HG3	1.99	0.44
1:A:287:LYS:HA	1:A:287:LYS:HD3	1.85	0.44
1:C:76:ILE:O	1:C:80:ILE:HG22	2.17	0.44
1:F:124:ASP:O	1:F:127:ILE:HG22	2.18	0.44
1:D:228:ASN:OD1	1:D:228:ASN:N	2.50	0.44
1:B:110:ILE:HD12	1:B:114:GLN:HE22	1.82	0.44
1:A:129:ILE:O	1:A:133:ILE:HG12	2.18	0.44
1:E:224:LEU:HD12	1:E:228:ASN:HD22	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:LYS:HA	1:D:36:LYS:HD3	1.80	0.43
1:B:123:VAL:O	1:B:127:ILE:HG22	2.19	0.43
1:B:171:LEU:HD23	1:B:171:LEU:HA	1.81	0.43
1:B:259:ASN:HB3	1:B:262:PHE:CE2	2.53	0.43
1:A:48:ARG:HG2	1:A:48:ARG:HH11	1.84	0.43
1:E:212:LEU:HD21	1:E:215:VAL:HG23	2.00	0.43
1:E:244:GLY:O	1:E:246:GLU:N	2.52	0.43
1:B:267:LEU:HD13	1:B:280:PHE:CE1	2.54	0.43
1:B:349:LEU:HD23	1:B:349:LEU:HA	1.87	0.43
1:D:249:LEU:O	1:D:297:LEU:HA	2.19	0.43
1:F:285:MET:HE2	1:F:285:MET:HB3	1.81	0.42
1:A:239:SER:O	1:A:241:GLN:NE2	2.52	0.42
1:B:72:HIS:O	1:B:76:ILE:HG12	2.19	0.42
1:B:361:ILE:N	1:B:361:ILE:HD12	2.34	0.42
1:E:117:PHE:CD2	1:E:121:GLY:HA3	2.54	0.42
1:B:35:LEU:HD21	1:B:127:ILE:HG13	2.02	0.42
1:F:79:LEU:HD12	1:F:79:LEU:O	2.19	0.42
1:D:70:HIS:HB3	1:D:116:HIS:CE1	2.54	0.42
1:D:167:ALA:O	1:D:170:THR:OG1	2.29	0.42
1:E:171:LEU:HD12	1:D:215:VAL:HG21	2.00	0.42
1:C:315:SER:O	1:C:315:SER:OG	2.33	0.42
1:A:89:VAL:HG22	1:A:133:ILE:HG22	2.02	0.42
1:C:267:LEU:HD11	1:C:278:ALA:HB3	2.02	0.42
1:B:145:ILE:HD12	1:B:208:PRO:HG3	2.01	0.42
1:F:172:TYR:HA	1:F:175:GLN:OE1	2.20	0.42
1:D:124:ASP:OD1	1:D:124:ASP:N	2.52	0.42
1:E:40:LEU:HD21	1:E:123:VAL:HG23	2.02	0.42
1:E:196:LEU:O	1:E:200:LEU:HB2	2.20	0.41
1:C:117:PHE:HB3	1:C:121:GLY:HA3	2.02	0.41
1:C:43:TRP:HB3	1:C:123:VAL:HG21	2.02	0.41
1:C:18:LEU:HD11	1:C:31:LEU:HD21	2.02	0.41
1:C:242:LEU:C	1:C:243:GLU:HG3	2.46	0.41
1:E:248:SER:O	1:E:249:LEU:HD23	2.21	0.41
1:E:297:LEU:HD12	1:E:297:LEU:HA	1.85	0.41
1:A:134:LEU:O	1:A:138:LEU:HD23	2.20	0.41
1:A:143:LYS:HE2	1:A:145:ILE:HD11	2.02	0.41
1:B:213:THR:OG1	1:B:234:LEU:O	2.38	0.41
1:B:387:LEU:HD12	1:B:387:LEU:HA	1.88	0.41
1:F:297:LEU:HD21	1:D:320:ARG:HD3	2.01	0.41
1:E:211:ASN:OD1	1:E:212:LEU:N	2.54	0.41
1:B:221:LYS:HA	1:B:221:LYS:HD2	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:174:LEU:HD23	1:E:174:LEU:HA	1.91	0.41
1:A:251:ILE:HG23	1:A:319:LEU:HD13	2.03	0.41
1:C:27:SER:HB2	1:C:30:GLU:HB2	2.02	0.41
1:F:271:MET:HE3	1:F:271:MET:HB3	1.99	0.41
1:E:62:VAL:HG11	1:E:127:ILE:HG21	2.03	0.41
1:F:217:PHE:HB3	1:F:229:LEU:HD12	2.01	0.41
1:F:231:ARG:HE	1:D:171:LEU:HD12	1.86	0.41
1:D:245:GLU:OE1	1:D:245:GLU:N	2.40	0.41
1:E:21:LEU:HD23	1:E:21:LEU:HA	1.86	0.41
1:A:223:GLU:OE1	1:A:223:GLU:N	2.44	0.41
1:F:13:SER:OG	1:F:16:MET:HG2	2.21	0.41
1:F:144:VAL:HG12	1:F:201:ALA:HA	2.02	0.41
1:E:112:ALA:O	1:E:115:VAL:HG12	2.20	0.41
1:E:194:ALA:O	1:E:198:THR:OG1	2.31	0.41
1:B:257:ASP:OD1	1:B:257:ASP:N	2.53	0.41
1:C:119:GLU:HG2	1:C:159:ARG:NH1	2.36	0.41
1:F:8:CYS:O	1:F:8:CYS:SG	2.79	0.41
1:A:78:GLY:HA2	1:A:81:ASN:HD21	1.86	0.40
1:B:149:LEU:HA	1:B:150:PRO:HD2	1.91	0.40
1:B:3:ILE:HD11	1:B:237:ILE:HD11	2.03	0.40
1:B:347:LEU:HD23	1:B:347:LEU:HA	1.91	0.40
1:F:218:GLY:N	1:F:230:LEU:O	2.53	0.40
1:D:41:GLU:OE1	1:D:42:GLY:N	2.55	0.40
1:D:218:GLY:N	1:D:230:LEU:O	2.49	0.40
1:A:102:ARG:O	1:A:102:ARG:NE	2.54	0.40
1:D:373:LYS:HE2	1:D:373:LYS:HB3	1.71	0.40
1:B:269:GLU:OE1	1:B:269:GLU:HA	2.22	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	378/414 (91%)	373 (99%)	5 (1%)	0	100	100
1	B	394/414 (95%)	389 (99%)	5 (1%)	0	100	100
1	C	398/414 (96%)	386 (97%)	12 (3%)	0	100	100
1	D	394/414 (95%)	384 (98%)	10 (2%)	0	100	100
1	E	396/414 (96%)	385 (97%)	8 (2%)	3 (1%)	16	50
1	F	379/414 (92%)	372 (98%)	7 (2%)	0	100	100
All	All	2339/2484 (94%)	2289 (98%)	47 (2%)	3 (0%)	49	80

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	247	SER
1	E	245	GLU
1	E	248	SER

5.3.2 Protein sidechains (i)

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	299/326 (92%)	298 (100%)	1 (0%)	86	85
1	B	310/326 (95%)	309 (100%)	1 (0%)	86	85
1	C	313/326 (96%)	313 (100%)	0	100	100
1	D	310/326 (95%)	308 (99%)	2 (1%)	78	80
1	E	313/326 (96%)	310 (99%)	3 (1%)	68	75
1	F	299/326 (92%)	299 (100%)	0	100	100
All	All	1844/1956 (94%)	1837 (100%)	7 (0%)	81	83

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	73	LEU
1	E	246	GLU

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Mol	Chain	Res	Type
1	E	248	SER
1	A	213	THR
1	B	328	VAL
1	D	108	HIS
1	D	299	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	118	HIS
1	E	228	ASN
1	A	81	ASN
1	A	114	GLN
1	B	28	GLN
1	B	114	GLN
1	B	370	GLN
1	C	81	ASN
1	C	118	HIS
1	F	70	HIS
1	D	52	GLN
1	D	68	GLN
1	D	70	HIS
1	D	99	ASN
1	D	330	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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 (i)

There are no chain breaks in this entry.

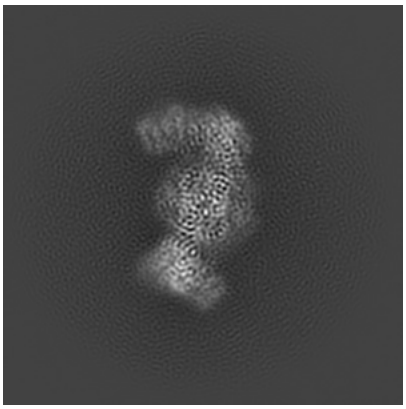
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-76280. These allow visual inspection of the internal detail of the map and identification of artifacts.

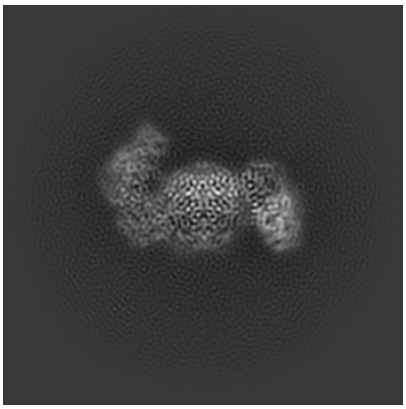
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

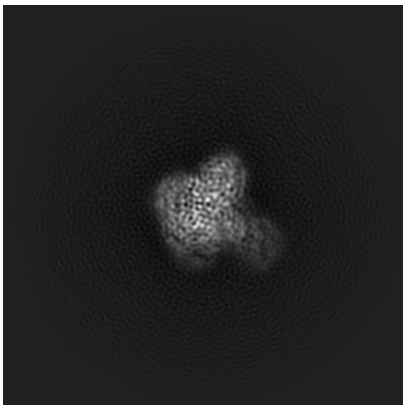
6.1.1 Primary map



X

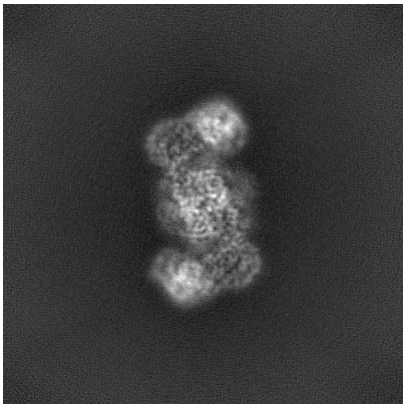


Y

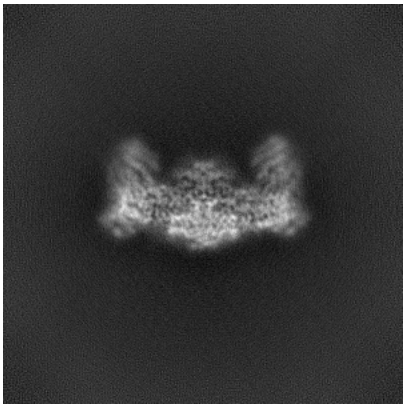


Z

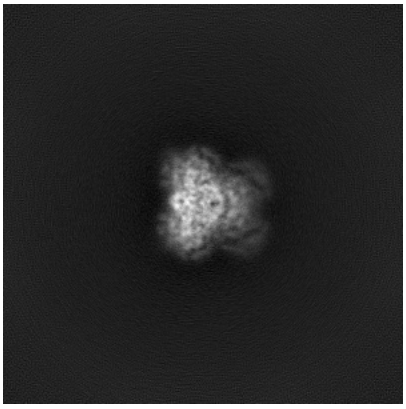
6.1.2 Raw map



X



Y

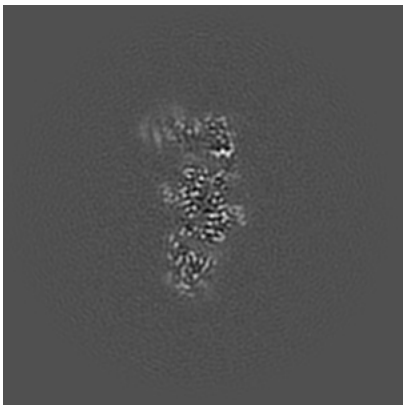


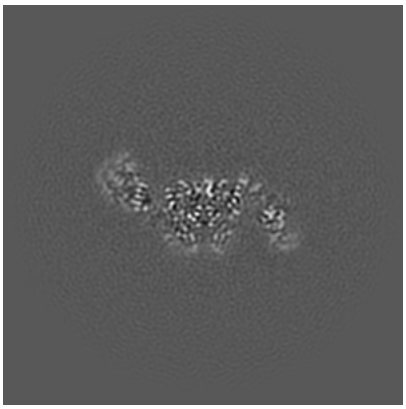
Z

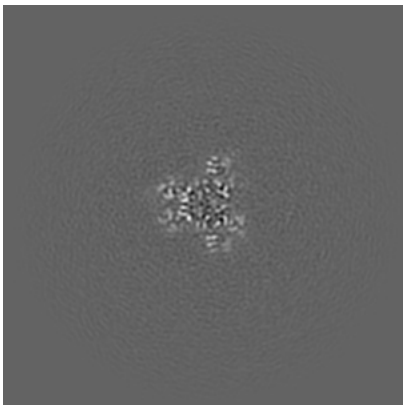
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

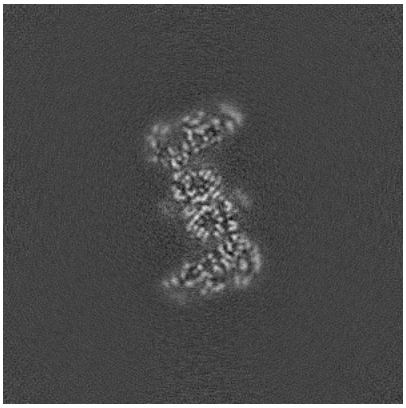
6.2.1 Primary map

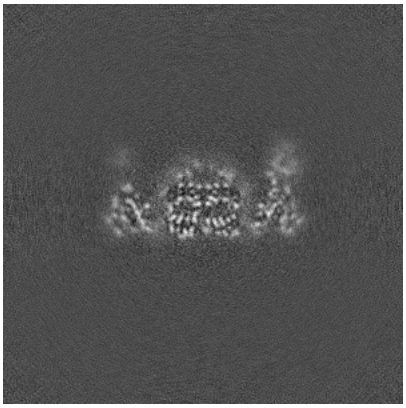

X Index: 192

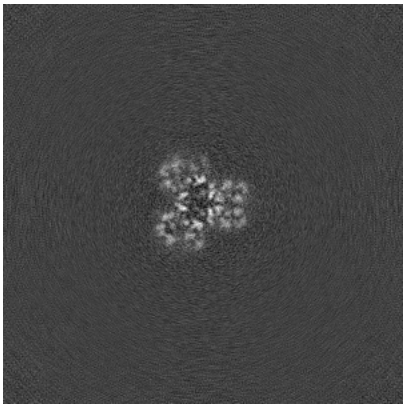

Y Index: 192


Z Index: 192

6.2.2 Raw map


X Index: 192

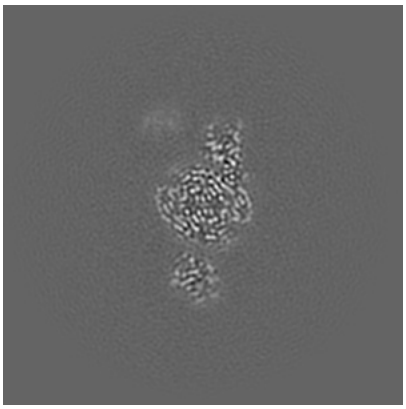

Y Index: 192

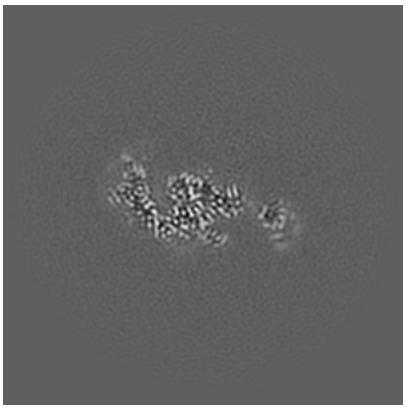

Z Index: 192

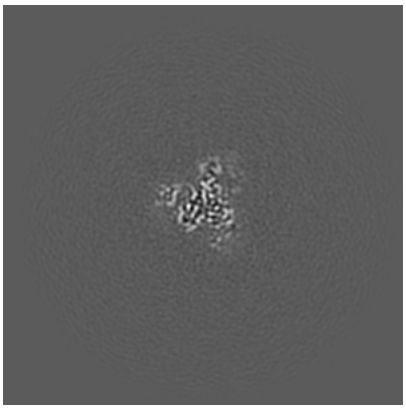
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

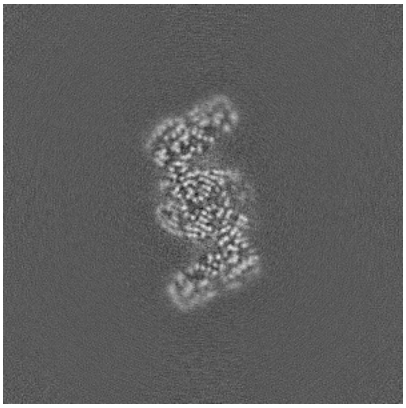
6.3.1 Primary map

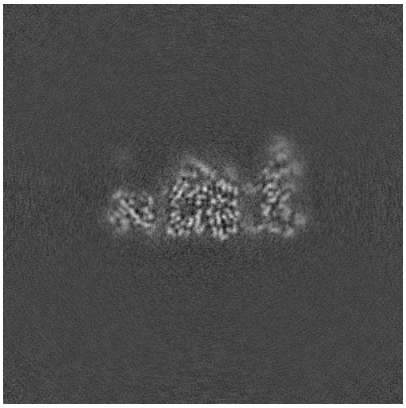

X Index: 203

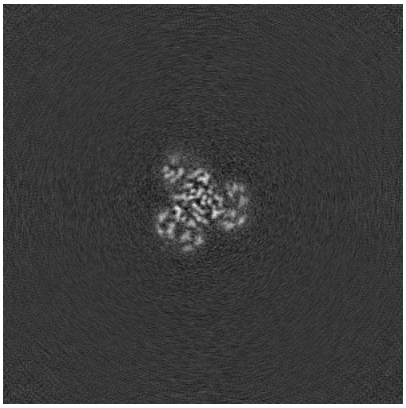

Y Index: 183


Z Index: 182

6.3.2 Raw map


X Index: 178

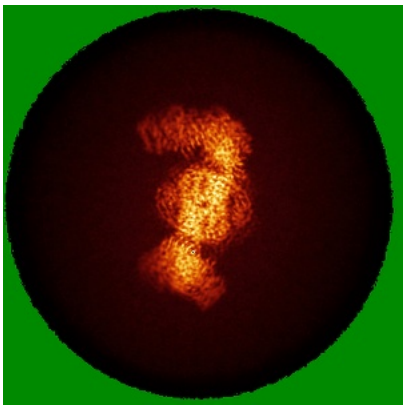

Y Index: 197

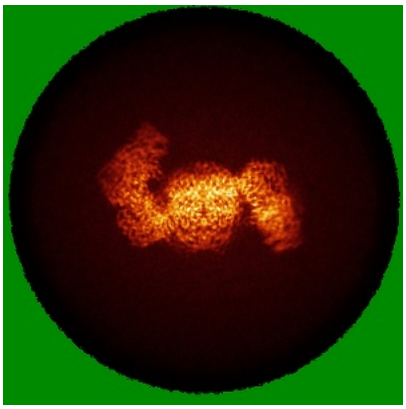

Z Index: 186

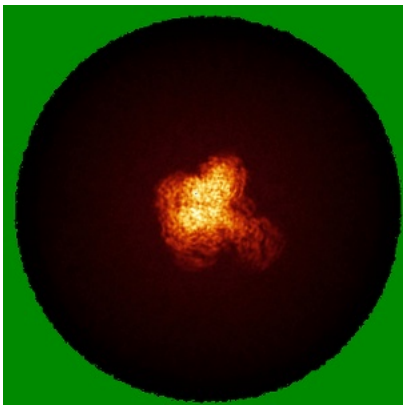
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

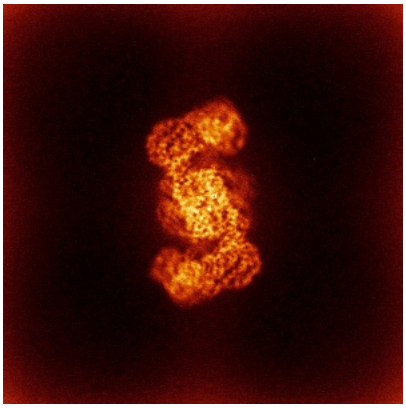
6.4.1 Primary map

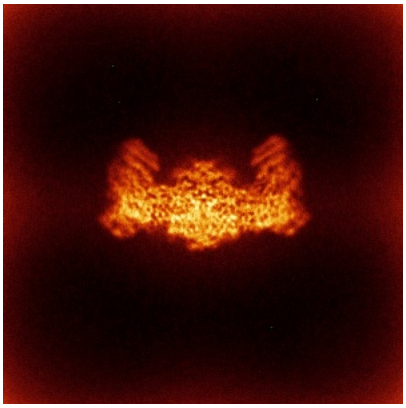

X

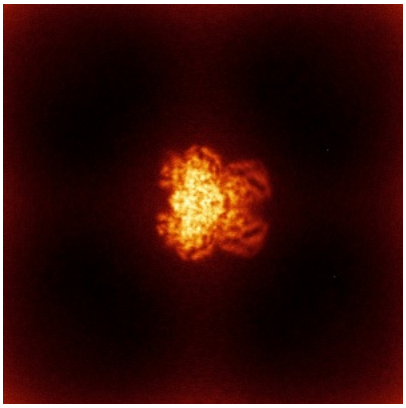

Y


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 0.12. 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.

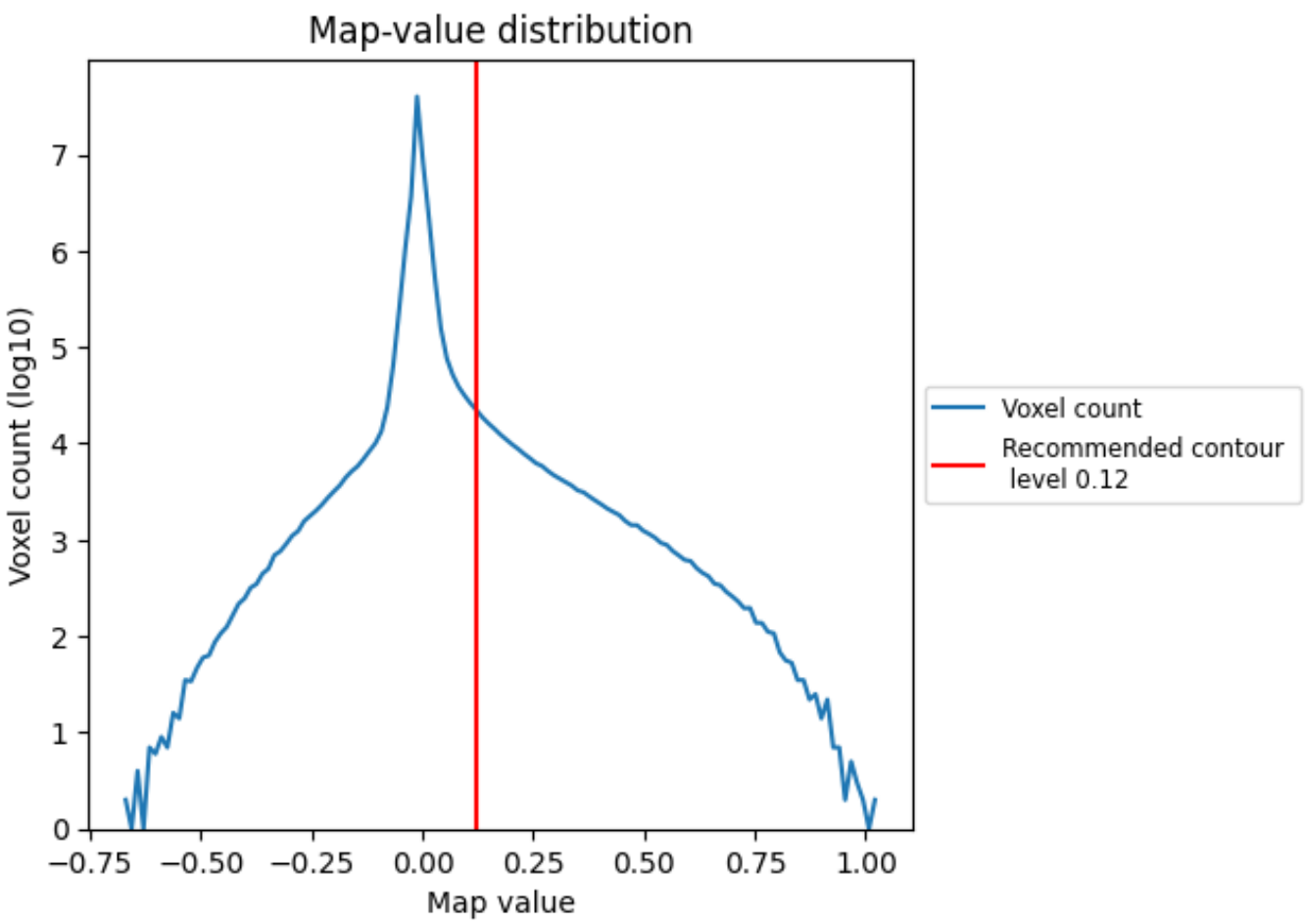
6.6 Mask visualisation (i)

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

7 Map analysis (i)

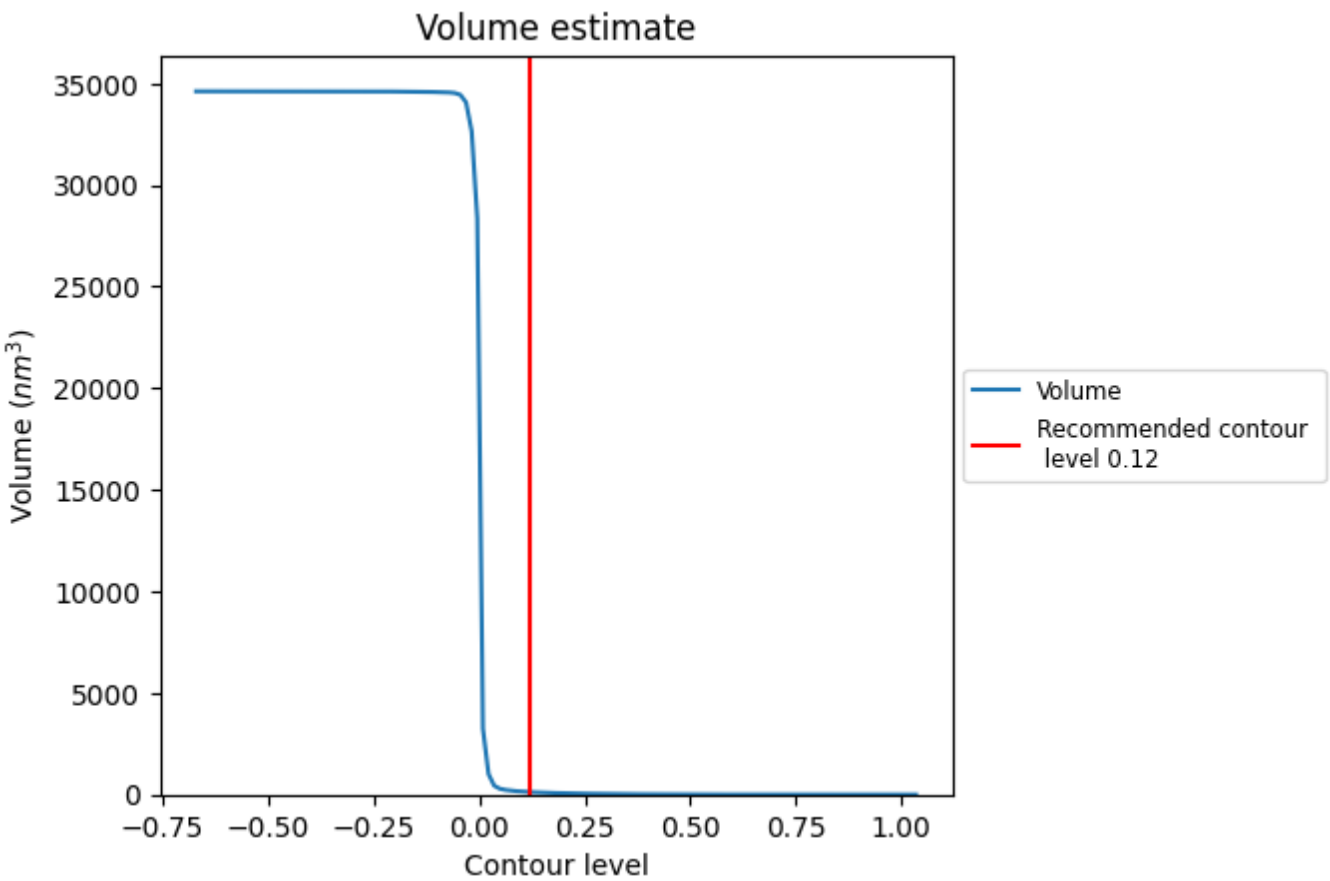
This section contains the results of statistical analysis of the map.

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.

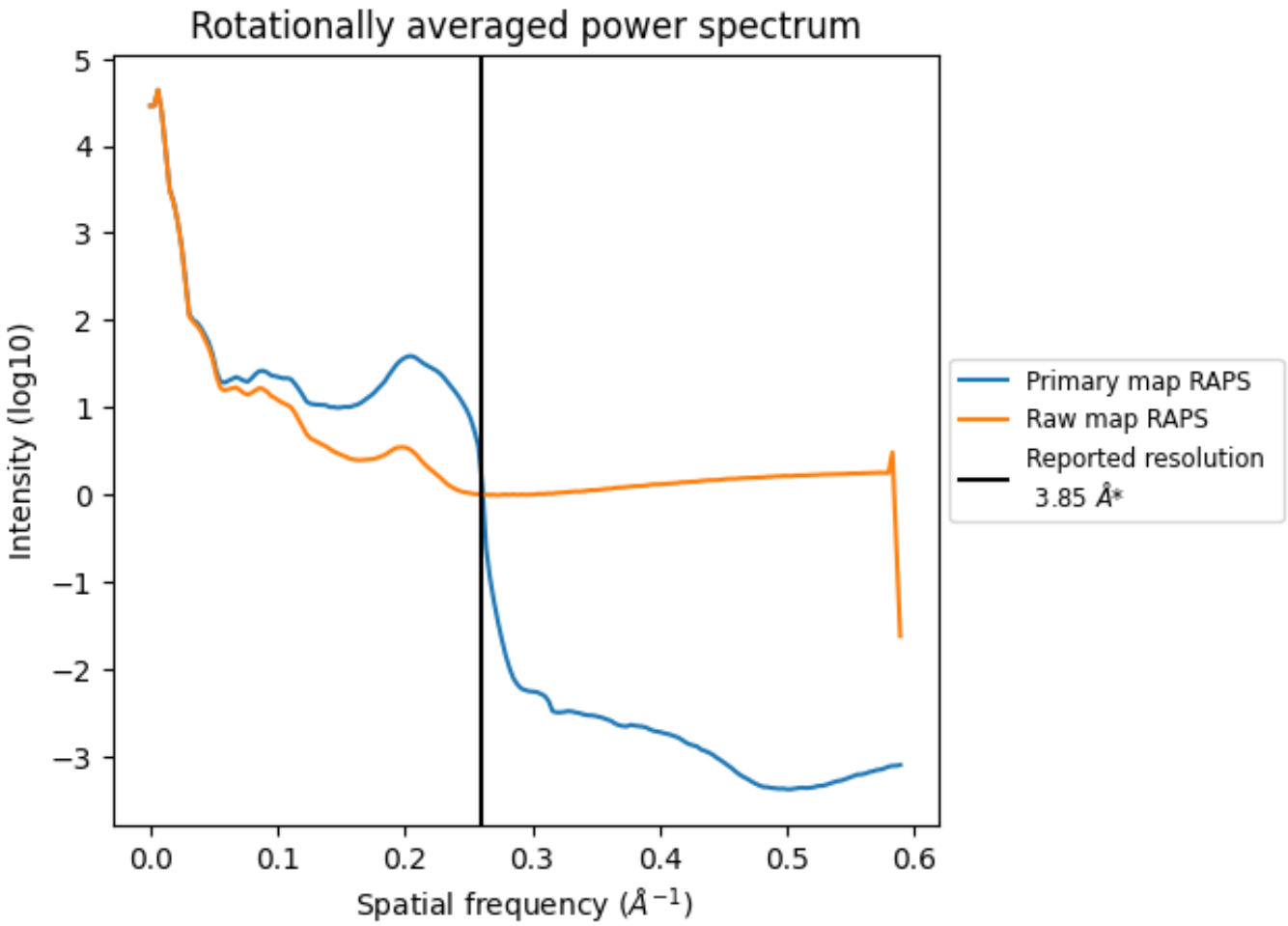
7.2 Volume estimate (i)



The volume at the recommended contour level is 123 nm^3; this corresponds to an approximate mass of 112 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)

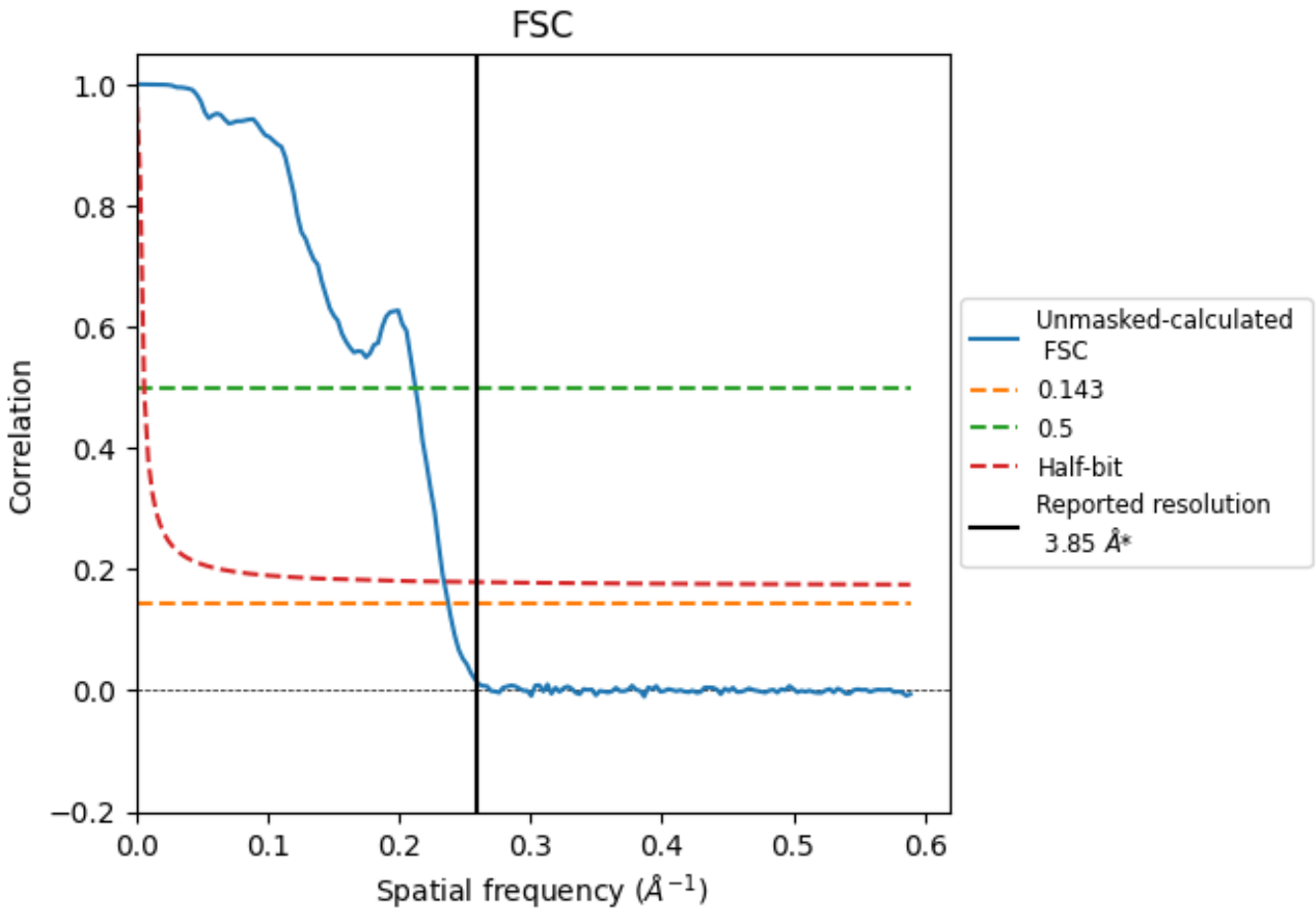


*Reported resolution corresponds to spatial frequency of 0.260 Å^{-1}

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 0.260 Å^{-1}

8.2 Resolution estimates (i)

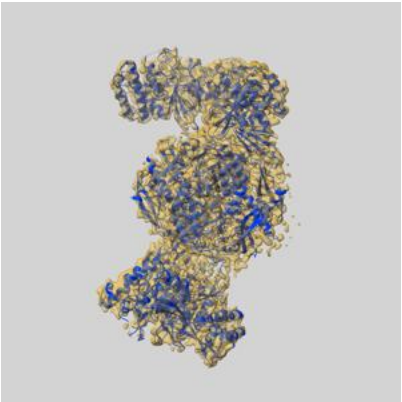
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.85	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.21	4.71	4.27

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

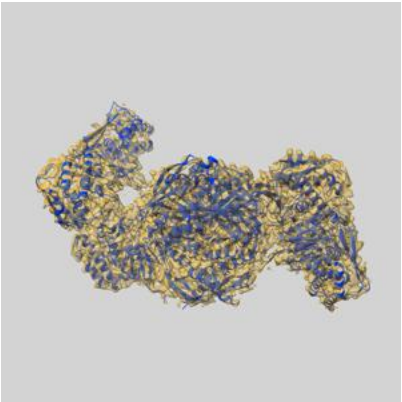
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-76280 and PDB model 12BJ. Per-residue inclusion information can be found in section 3 on page 8.

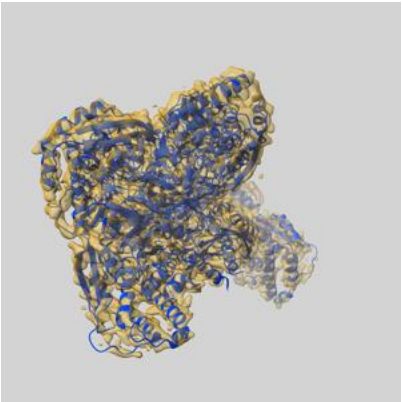
9.1 Map-model overlay (i)



X



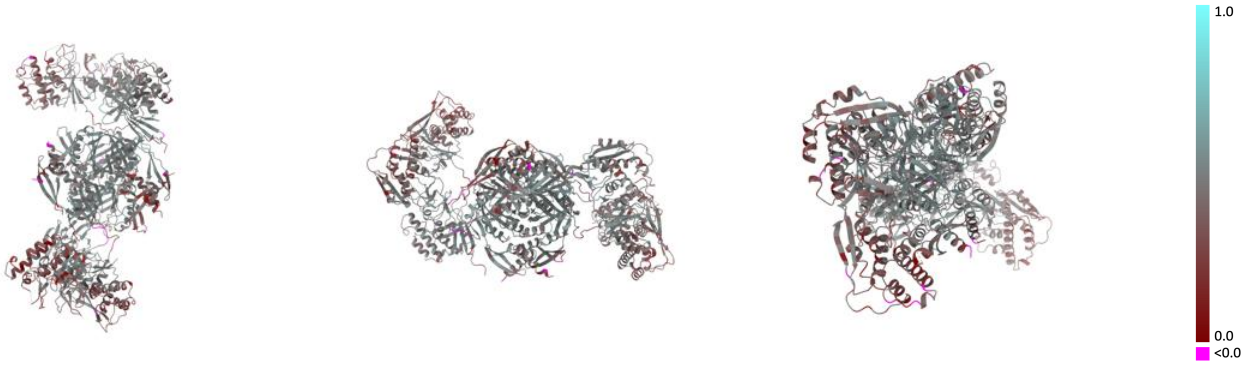
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.12 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.

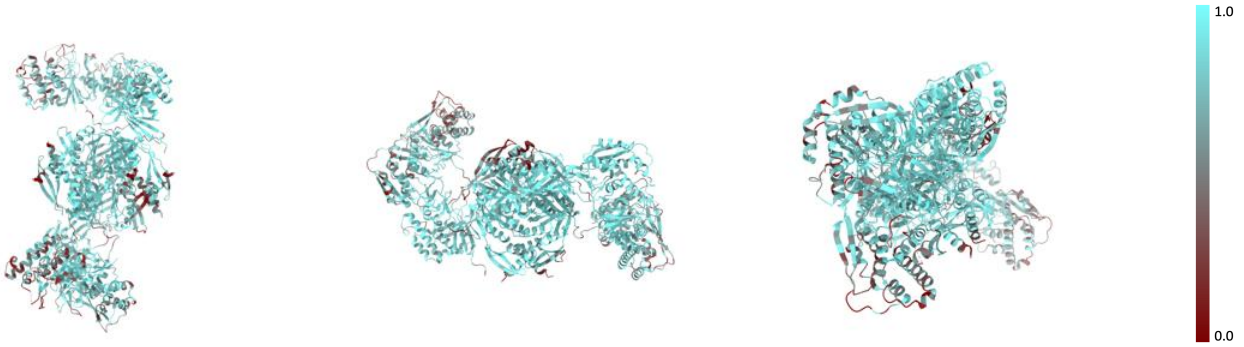
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

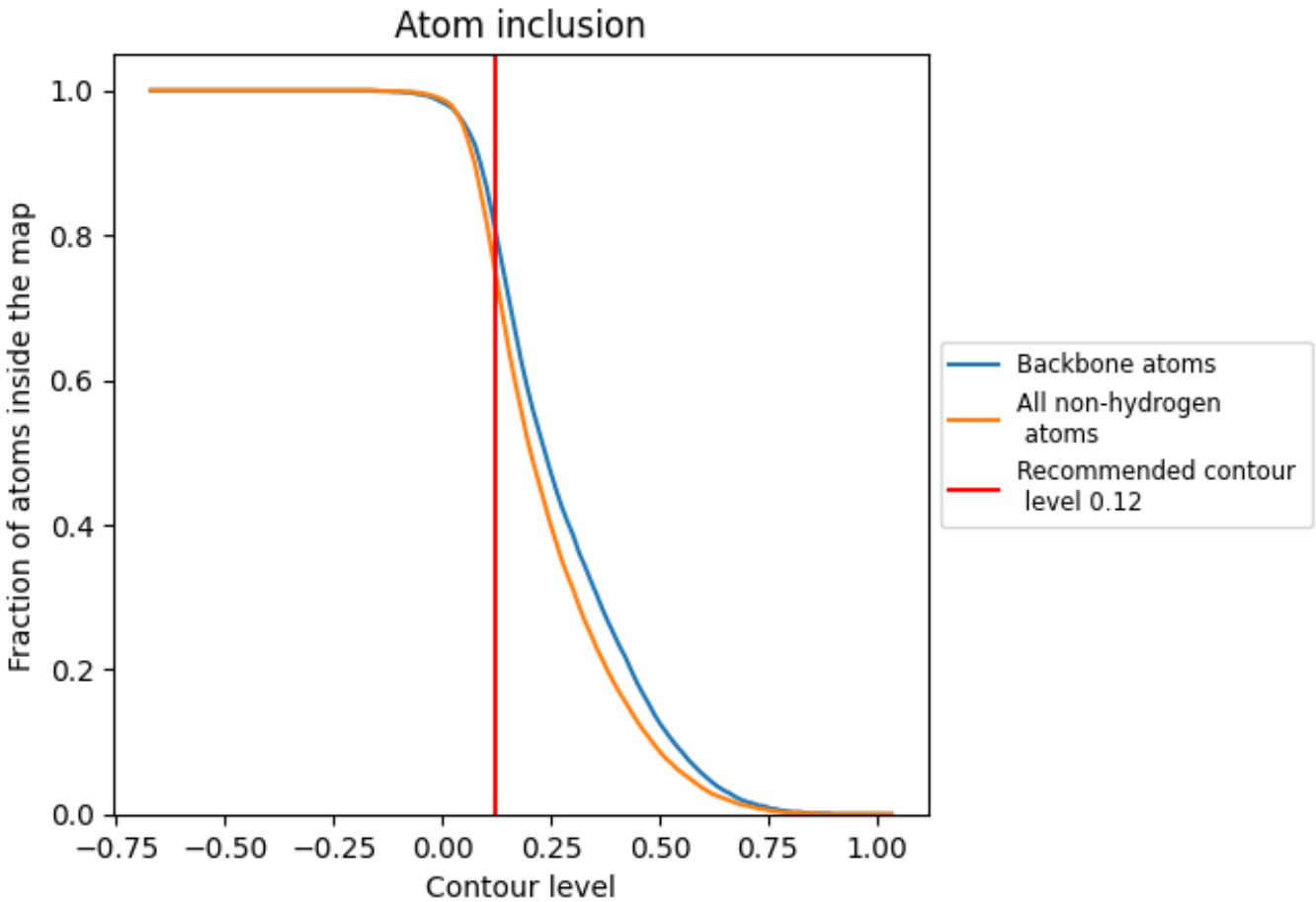
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.12).

9.4 Atom inclusion (i)



At the recommended contour level, 82% of all backbone atoms, 76% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.12) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7570	0.4360
A	0.7090	0.4140
B	0.7640	0.4380
C	0.8360	0.4790
D	0.7150	0.4140
E	0.8210	0.4670
F	0.6890	0.4040