Full wwPDB EM Validation Report ⓘ

Oct 21, 2026 – 02:26 PM EDT

PDB ID	:	10GF / pdb_000010gf
EMDB ID	:	EMD-75149
Title	:	The eukaryotic translation initiation factor 2B from Homo sapiens with a L179V mutation in the delta subunit, I-state
Authors	:	Dalwadi, U.; Boone, M.; Croll, T.I.; Laughlin, T.G.; Frost, A.; Walter, P.
Deposited on	:	2026-01-19
Resolution	:	2.50 Å(reported)
Based on initial model	:	7L70

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:

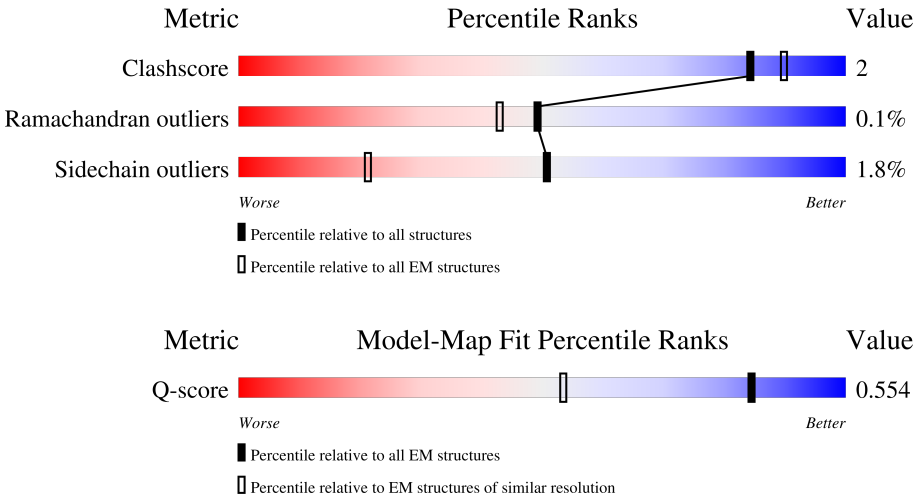
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4-5-2 with Phenix2.2
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

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 2.50 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	-
Q-score	-	25397	7115 (2.00 - 3.00)

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 <=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	721	54% 6% 39%
1	B	721	55% 5% 39%
2	E	523	59% 6% 34%
2	F	523	7% 59% 6% 34%

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Mol	Chain	Length	Quality of chain
3	G	322	15% 88% 7% 5%
3	H	322	14% 84% 11% 5%
4	I	452	29% 81% 6% 13%
4	J	452	53% 85% 13%
5	C	368	80% 7% 12%
5	D	368	78% 10% 12%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	PO4	D	1000	-	-	X	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 28056 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 Translation initiation factor eIF-2B subunit epsilon.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	437	Total 3423	C 2167	N 606	O 635	S 15	0	0
1	B	437	Total 3423	C 2167	N 606	O 635	S 15	0	0

- Molecule 2 is a protein called Translation initiation factor eIF2B subunit delta.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	345	Total 2682	C 1695	N 477	O 496	S 14	0	0
2	F	345	Total 2682	C 1695	N 477	O 496	S 14	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	179	VAL	LEU	engineered mutation	UNP Q9UI10
F	179	VAL	LEU	engineered mutation	UNP Q9UI10

- Molecule 3 is a protein called Translation initiation factor eIF2B subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	305	Total 2368	C 1517	N 393	O 446	S 12	0	0
3	H	305	Total 2368	C 1517	N 393	O 446	S 12	0	0

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	-16	MET	-	initiating methionine	UNP Q14232
G	-15	HIS	-	expression tag	UNP Q14232
G	-14	HIS	-	expression tag	UNP Q14232

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-13	HIS	-	expression tag	UNP Q14232
G	-12	HIS	-	expression tag	UNP Q14232
G	-11	HIS	-	expression tag	UNP Q14232
G	-10	HIS	-	expression tag	UNP Q14232
G	-9	GLY	-	expression tag	UNP Q14232
G	-8	GLY	-	expression tag	UNP Q14232
G	-7	GLY	-	expression tag	UNP Q14232
G	-6	SER	-	expression tag	UNP Q14232
G	-5	GLU	-	expression tag	UNP Q14232
G	-4	ASN	-	expression tag	UNP Q14232
G	-3	LEU	-	expression tag	UNP Q14232
G	-2	TYR	-	expression tag	UNP Q14232
G	-1	PHE	-	expression tag	UNP Q14232
G	0	GLN	-	expression tag	UNP Q14232
H	-16	MET	-	initiating methionine	UNP Q14232
H	-15	HIS	-	expression tag	UNP Q14232
H	-14	HIS	-	expression tag	UNP Q14232
H	-13	HIS	-	expression tag	UNP Q14232
H	-12	HIS	-	expression tag	UNP Q14232
H	-11	HIS	-	expression tag	UNP Q14232
H	-10	HIS	-	expression tag	UNP Q14232
H	-9	GLY	-	expression tag	UNP Q14232
H	-8	GLY	-	expression tag	UNP Q14232
H	-7	GLY	-	expression tag	UNP Q14232
H	-6	SER	-	expression tag	UNP Q14232
H	-5	GLU	-	expression tag	UNP Q14232
H	-4	ASN	-	expression tag	UNP Q14232
H	-3	LEU	-	expression tag	UNP Q14232
H	-2	TYR	-	expression tag	UNP Q14232
H	-1	PHE	-	expression tag	UNP Q14232
H	0	GLN	-	expression tag	UNP Q14232

- Molecule 4 is a protein called Translation initiation factor eIF-2B subunit gamma.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	I	391	Total	C	N	O	S	0	0
			3026	1920	516	566	24		
4	J	391	Total	C	N	O	S	0	0
			3026	1920	516	566	24		

- Molecule 5 is a protein called Translation initiation factor eIF-2B subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	C	322	Total	C	N	O	S	0	0
			2523	1596	442	470	15		
5	D	322	Total	C	N	O	S	0	0
			2523	1596	442	470	15		

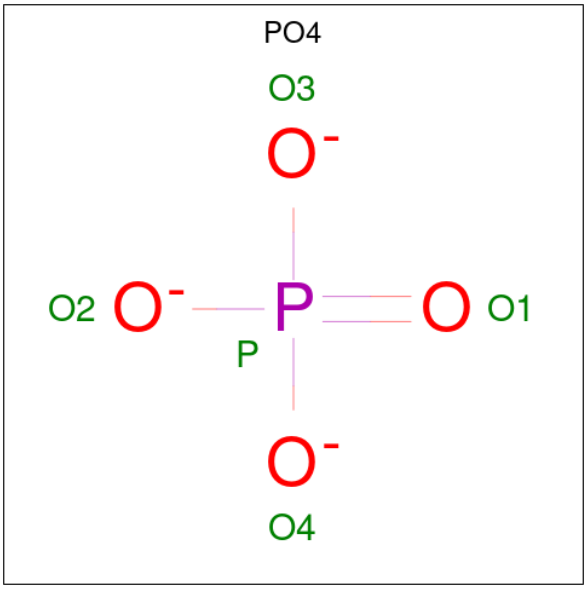
There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-16	MET	-	initiating methionine	UNP P49770
C	-15	HIS	-	expression tag	UNP P49770
C	-14	HIS	-	expression tag	UNP P49770
C	-13	HIS	-	expression tag	UNP P49770
C	-12	HIS	-	expression tag	UNP P49770
C	-11	HIS	-	expression tag	UNP P49770
C	-10	HIS	-	expression tag	UNP P49770
C	-9	GLY	-	expression tag	UNP P49770
C	-8	GLY	-	expression tag	UNP P49770
C	-7	GLY	-	expression tag	UNP P49770
C	-6	SER	-	expression tag	UNP P49770
C	-5	GLU	-	expression tag	UNP P49770
C	-4	ASN	-	expression tag	UNP P49770
C	-3	LEU	-	expression tag	UNP P49770
C	-2	TYR	-	expression tag	UNP P49770
C	-1	PHE	-	expression tag	UNP P49770
C	0	GLN	-	expression tag	UNP P49770
C	1	SER	-	expression tag	UNP P49770
D	-16	MET	-	initiating methionine	UNP P49770
D	-15	HIS	-	expression tag	UNP P49770
D	-14	HIS	-	expression tag	UNP P49770
D	-13	HIS	-	expression tag	UNP P49770
D	-12	HIS	-	expression tag	UNP P49770
D	-11	HIS	-	expression tag	UNP P49770
D	-10	HIS	-	expression tag	UNP P49770
D	-9	GLY	-	expression tag	UNP P49770
D	-8	GLY	-	expression tag	UNP P49770
D	-7	GLY	-	expression tag	UNP P49770
D	-6	SER	-	expression tag	UNP P49770
D	-5	GLU	-	expression tag	UNP P49770
D	-4	ASN	-	expression tag	UNP P49770
D	-3	LEU	-	expression tag	UNP P49770
D	-2	TYR	-	expression tag	UNP P49770
D	-1	PHE	-	expression tag	UNP P49770
D	0	GLN	-	expression tag	UNP P49770
D	1	SER	-	expression tag	UNP P49770

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	AltConf
6	E	1	Total Zn 1 1	0
6	F	1	Total Zn 1 1	0

- Molecule 7 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	AltConf
7	C	1	Total O P 5 4 1	0
7	D	1	Total O P 5 4 1	0

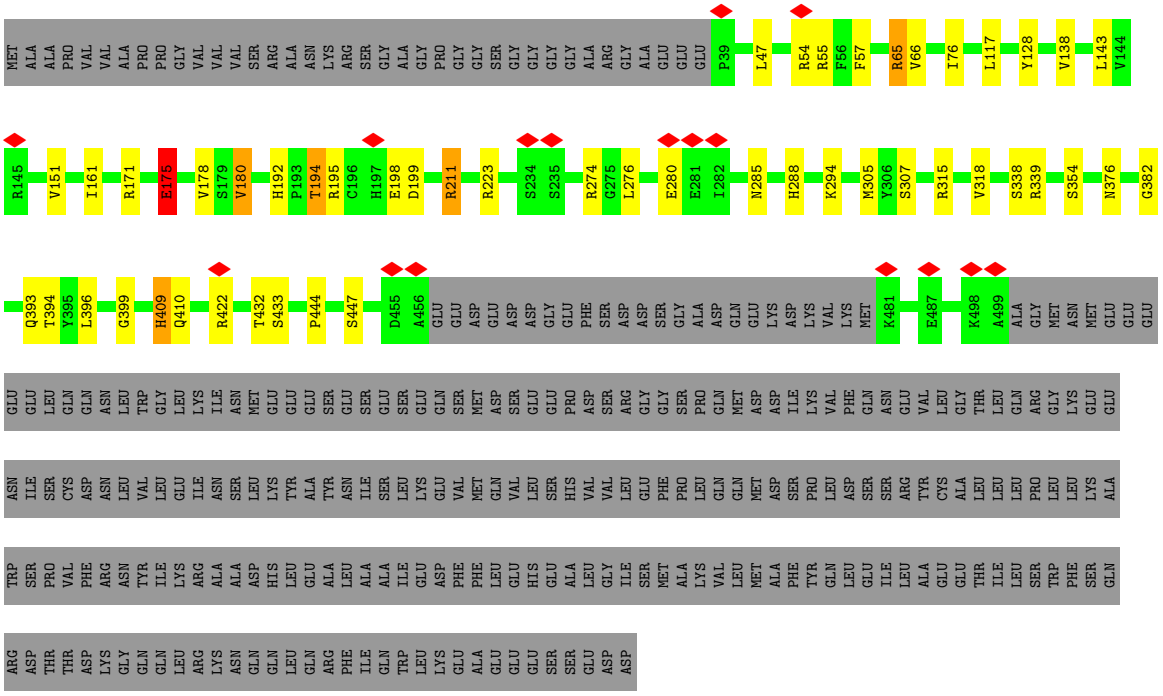
3 Residue-property plots (i)

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: Translation initiation factor eIF-2B subunit epsilon

Chain A:

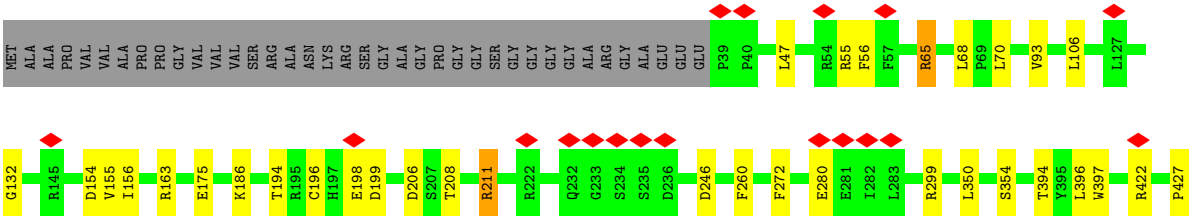




• Molecule 1: Translation initiation factor eIF-2B subunit epsilon

Chain B:



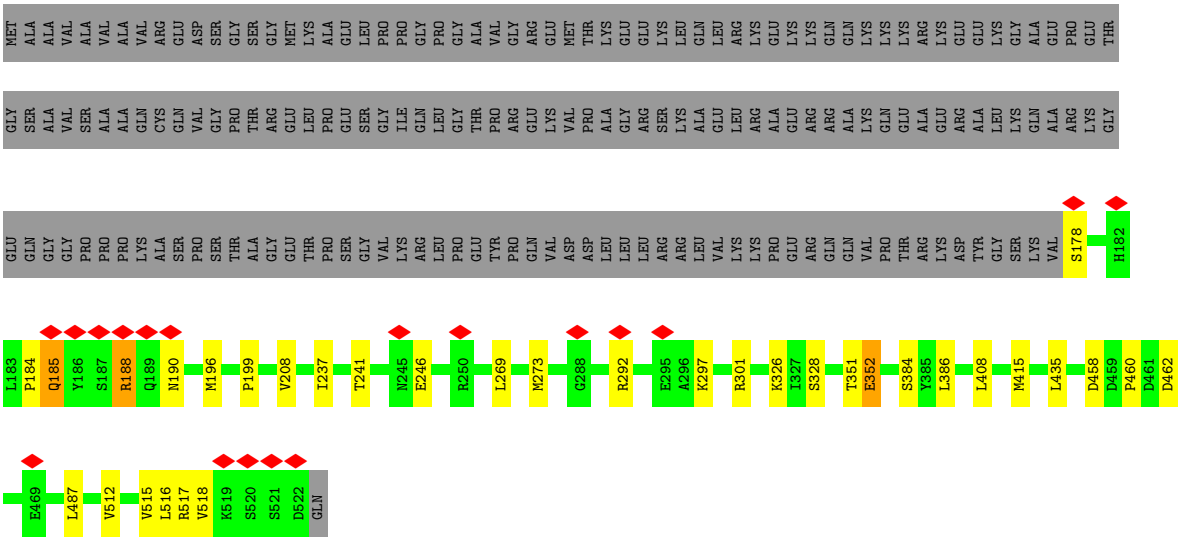




• Molecule 2: Translation initiation factor eIF2B subunit delta

Chain E:

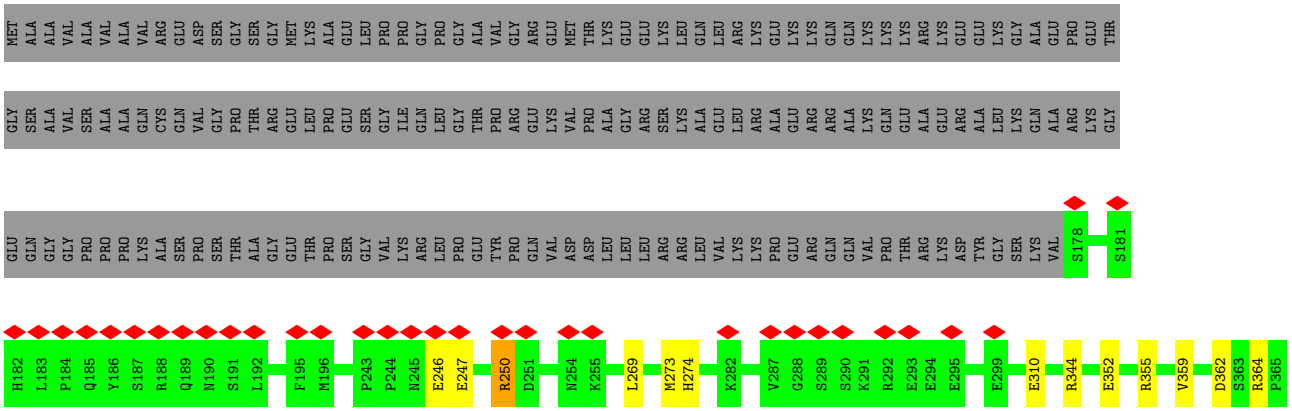


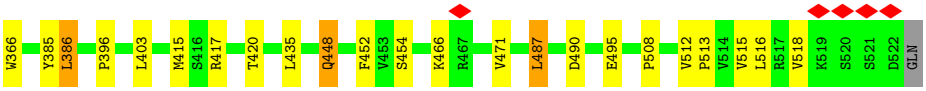


• Molecule 2: Translation initiation factor eIF2B subunit delta

Chain F:



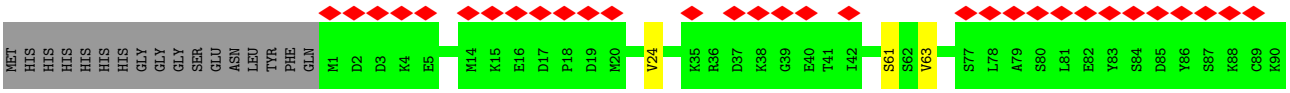




• Molecule 3: Translation initiation factor eIF2B subunit alpha

Chain G:



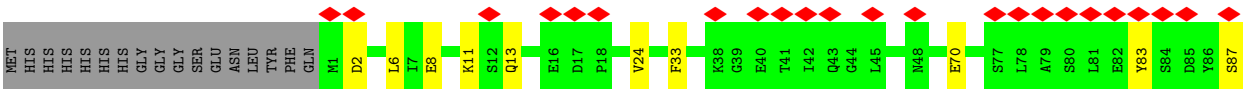


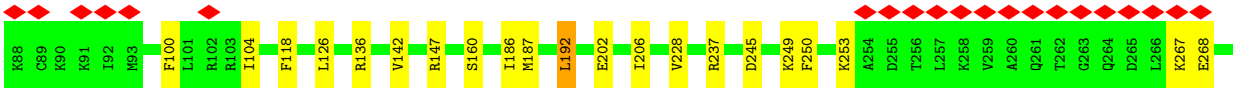


• Molecule 3: Translation initiation factor eIF2B subunit alpha

Chain H:





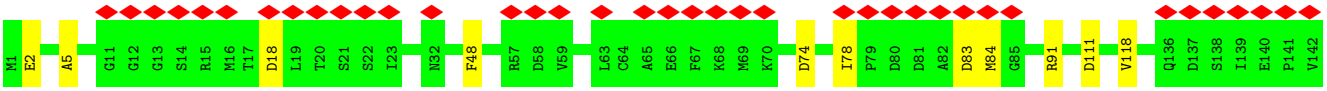


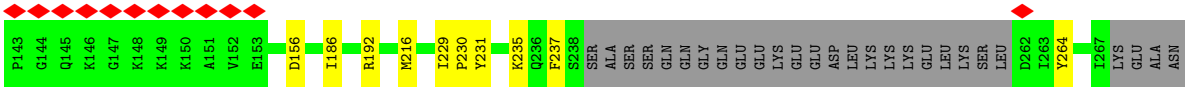


• Molecule 4: Translation initiation factor eIF-2B subunit gamma

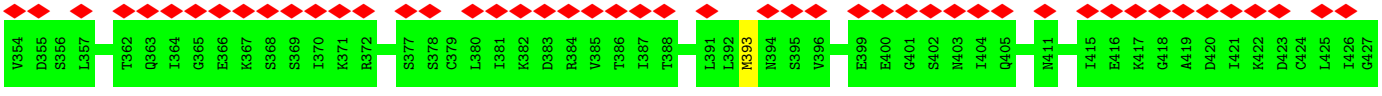
Chain I:

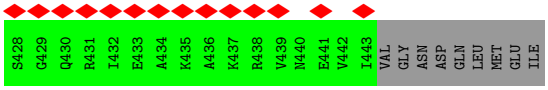






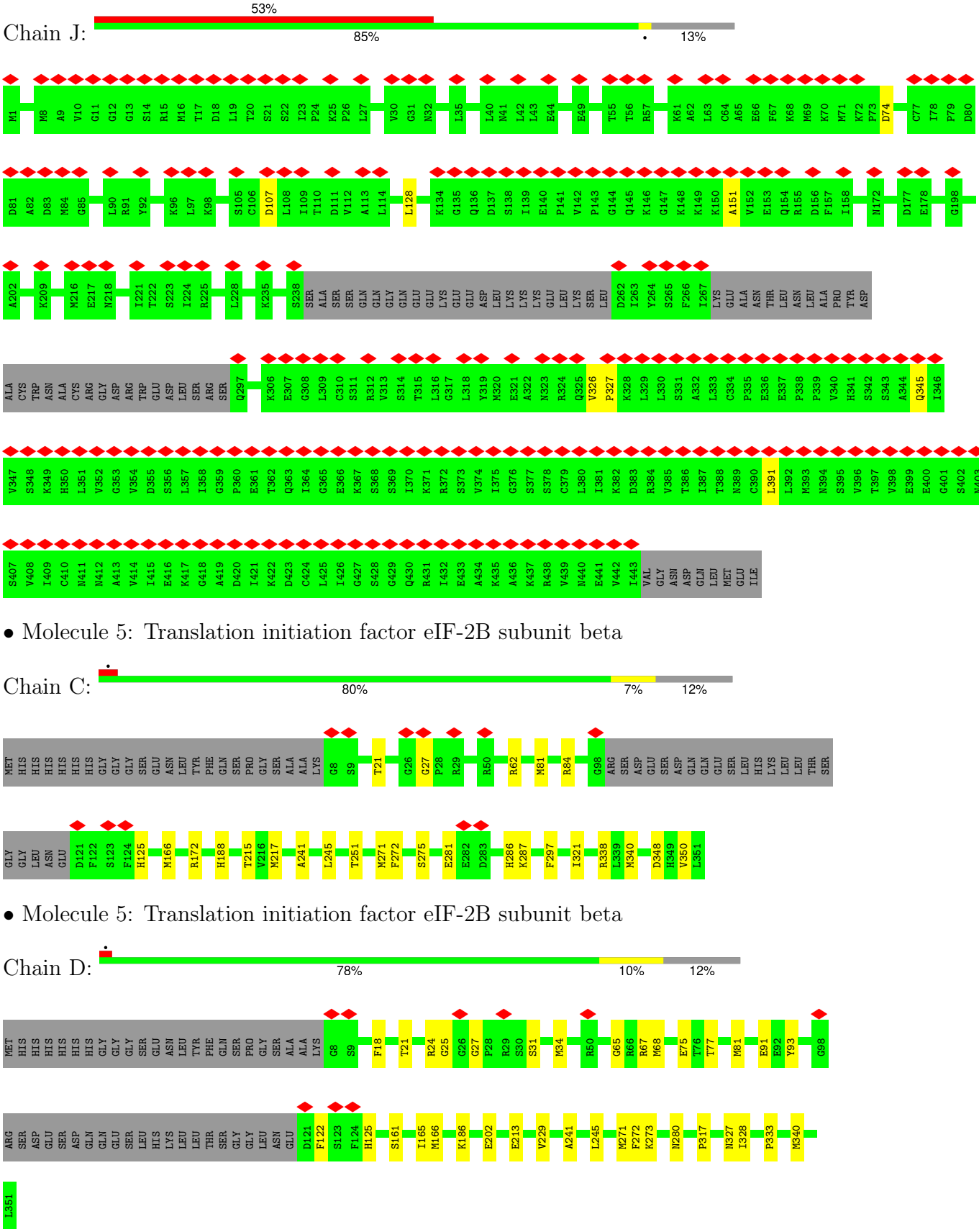






• Molecule 4: Translation initiation factor eIF-2B subunit gamma

Chain J:



• Molecule 5: Translation initiation factor eIF-2B subunit beta

Chain C:

• Molecule 5: Translation initiation factor eIF-2B subunit beta

Chain D:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	60754	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 ($e^-/Å^2$)	46.7	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1600	Depositor
Magnification	135000	Depositor
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	0.935	Depositor
Minimum map value	-0.476	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.030	Depositor
Recommended contour level	0.165	Depositor
Map size (Å)	371.44, 371.44, 371.44	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.9286, 0.9286, 0.9286	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, ZN

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	0.74	1/3497 (0.0%)	1.18	8/4758 (0.2%)
1	B	0.64	0/3497	1.12	3/4758 (0.1%)
2	E	0.76	1/2732 (0.0%)	1.15	2/3714 (0.1%)
2	F	0.72	0/2732	1.13	6/3714 (0.2%)
3	G	0.72	0/2405	1.18	3/3247 (0.1%)
3	H	0.72	1/2405 (0.0%)	1.17	5/3247 (0.2%)
4	I	0.59	0/3069	1.05	5/4140 (0.1%)
4	J	0.55	0/3069	0.96	2/4140 (0.0%)
5	C	0.77	1/2572 (0.0%)	1.19	4/3477 (0.1%)
5	D	0.78	0/2572	1.22	3/3477 (0.1%)
All	All	0.70	4/28550 (0.0%)	1.13	41/38672 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	6
1	B	0	4
2	E	0	3
2	F	0	4
3	G	0	1
4	I	0	1
5	C	0	2
5	D	0	4
All	All	0	25

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	307	SER	CA-CB	-5.18	1.44	1.53
3	H	160	SER	CA-CB	-5.16	1.44	1.53
2	E	384	SER	CA-CB	-5.13	1.45	1.54
5	C	275	SER	CA-CB	-5.04	1.45	1.53

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	194	THR	OG1-CB-CG2	-9.91	89.47	109.30
2	E	352	GLU	N-CA-CB	7.06	122.43	110.49
5	D	77	THR	CA-CB-OG1	-6.72	99.53	109.60
3	H	245	ASP	CA-CB-CG	6.62	119.22	112.60
2	F	352	GLU	N-CA-CB	6.35	121.41	114.17
1	B	194	THR	OG1-CB-CG2	-6.32	96.65	109.30
2	E	188	ARG	NE-CZ-NH2	6.03	124.62	119.20
2	F	448	GLN	N-CA-CB	-5.96	100.34	111.13
5	C	84	ARG	NE-CZ-NH2	5.91	124.52	119.20
3	H	2	ASP	CA-CB-CG	5.88	118.48	112.60
1	B	246	ASP	CA-CB-CG	5.85	118.45	112.60
3	H	136	ARG	CA-CB-CG	-5.80	102.50	114.10
5	C	281	GLU	N-CA-CB	5.77	118.87	109.69
3	H	202	GLU	CB-CG-CD	5.75	122.38	112.60
3	H	274	ASP	CA-CB-CG	5.71	118.31	112.60
3	G	302	LYS	CB-CA-C	-5.67	101.39	110.79
3	G	245	ASP	CA-CB-CG	5.58	118.18	112.60
4	I	83	ASP	CA-CB-CG	5.57	118.17	112.60
4	I	192	ARG	N-CA-CB	-5.56	102.43	112.27
2	F	487	LEU	N-CA-CB	-5.55	101.45	109.83
1	B	154	ASP	CA-CB-CG	5.54	118.14	112.60
1	A	180	VAL	N-CA-CB	-5.52	102.43	111.98
1	A	128	TYR	CB-CA-C	-5.52	101.29	109.90
5	D	317	PRO	CB-CA-C	-5.50	106.22	111.39
2	F	310	GLU	CB-CA-C	-5.45	102.50	110.95
2	F	490	ASP	CA-CB-CG	5.44	118.04	112.60
1	A	175	GLU	CB-CA-C	5.39	119.27	111.82
4	I	74	ASP	CA-CB-CG	5.39	118.00	112.60
1	A	195	ARG	CA-CB-CG	-5.33	103.43	114.10
4	J	74	ASP	CA-CB-CG	5.32	117.92	112.60
1	A	199	ASP	CA-CB-CG	5.31	117.91	112.60
5	C	251	THR	CA-CB-OG1	-5.29	101.67	109.60
3	G	110	LYS	N-CA-CB	5.28	117.72	110.07
1	A	409	HIS	CA-CB-CG	-5.25	108.55	113.80
4	I	156	ASP	CA-CB-CG	5.23	117.83	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	I	111	ASP	CA-CB-CG	5.20	117.80	112.60
5	C	287	LYS	N-CA-C	-5.10	101.65	109.76
2	F	508	PRO	CB-CA-C	5.07	117.62	110.98
5	D	186	LYS	N-CA-CB	5.06	118.54	110.65
1	A	294	LYS	N-CA-CB	-5.05	103.11	111.49
4	J	107	ASP	CA-CB-CG	5.03	117.63	112.60

There are no chirality outliers.

All (25) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	175	GLU	Peptide
1	A	211	ARG	Sidechain
1	A	274	ARG	Sidechain
1	A	315	ARG	Sidechain
1	A	54	ARG	Sidechain
1	A	65	ARG	Sidechain
1	B	163	ARG	Sidechain
1	B	211	ARG	Sidechain
1	B	299	ARG	Sidechain
1	B	65	ARG	Sidechain
5	C	172	ARG	Sidechain
5	C	338	ARG	Sidechain
5	D	24	ARG	Sidechain,Peptide
5	D	245	LEU	Peptide
5	D	67	ARG	Sidechain
2	E	292	ARG	Sidechain
2	E	301	ARG	Sidechain
2	E	517	ARG	Sidechain
2	F	250	ARG	Sidechain
2	F	344	ARG	Sidechain
2	F	355	ARG	Sidechain
2	F	364	ARG	Sidechain
3	G	136	ARG	Sidechain
4	I	299	ARG	Sidechain

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	3423	0	3419	21	0
1	B	3423	0	3419	19	0
2	E	2682	0	2735	13	0
2	F	2682	0	2735	13	0
3	G	2368	0	2447	12	0
3	H	2368	0	2447	15	0
4	I	3026	0	3144	9	0
4	J	3026	0	3144	2	0
5	C	2523	0	2527	15	0
5	D	2523	0	2527	17	0
6	E	1	0	0	0	0
6	F	1	0	0	0	0
7	C	5	0	0	0	0
7	D	5	0	0	2	0
All	All	28056	0	28544	128	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 2.

All (128) 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)
2:E:188:ARG:NH1	2:E:462:ASP:OD2	2.15	0.78
1:A:171:ARG:HD2	1:A:288:HIS:CD2	2.33	0.64
5:D:241:ALA:HA	5:D:340:MET:HE3	1.81	0.63
2:E:515:VAL:HG12	2:E:516:LEU:HD23	1.82	0.61
1:B:394:THR:HG22	1:B:396:LEU:CD1	2.30	0.61
1:A:211:ARG:HH11	1:A:280:GLU:CD	2.10	0.60
1:A:138:VAL:HA	1:A:143:LEU:HD12	1.84	0.59
1:B:394:THR:HG22	1:B:396:LEU:HD12	1.87	0.56
5:C:241:ALA:HA	5:C:340:MET:HE3	1.87	0.56
1:A:47:LEU:C	1:A:47:LEU:HD23	2.31	0.56
5:D:21:THR:HG22	5:D:27:GLY:HA2	1.88	0.55
5:D:166:MET:HE3	5:D:229:VAL:HG21	1.88	0.55
1:B:186:LYS:HE3	5:C:297:PHE:CD2	2.41	0.55
5:C:166:MET:HE2	5:C:217:MET:HE1	1.89	0.55
3:H:285:PHE:CD2	3:H:290:VAL:HG22	2.41	0.55
2:E:178:SER:N	5:D:213:GLU:OE2	2.40	0.55
5:D:65:GLY:HA2	5:D:68:MET:HE3	1.89	0.55
2:F:359:VAL:CG1	2:F:386:LEU:HD21	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:126:LEU:HD13	3:G:186:ILE:CD1	2.36	0.55
5:D:25:GLY:O	5:D:27:GLY:N	2.38	0.54
3:G:126:LEU:HD13	3:G:186:ILE:HD12	1.88	0.54
5:D:21:THR:CG2	5:D:27:GLY:HA2	2.40	0.52
5:C:81:MET:HE3	5:C:272:PHE:HB2	1.92	0.52
3:G:192:LEU:HD23	3:G:192:LEU:N	2.23	0.52
1:B:55:ARG:NH1	1:B:198:GLU:OE2	2.44	0.50
3:H:126:LEU:HD13	3:H:186:ILE:CD1	2.42	0.50
4:I:91:ARG:HA	4:I:216:MET:HE1	1.93	0.50
5:C:166:MET:HE2	5:C:217:MET:CE	2.42	0.50
1:A:338:SER:O	1:A:339:ARG:C	2.55	0.50
1:B:211:ARG:HH11	1:B:280:GLU:CD	2.20	0.49
5:D:18:PHE:CE2	5:D:68:MET:SD	3.06	0.49
3:G:61:SER:HB2	3:G:249:LYS:O	2.12	0.49
2:F:452:PHE:CE2	2:F:454:SER:HB3	2.47	0.49
2:F:247:GLU:OE1	2:F:250:ARG:NH2	2.46	0.49
3:H:70:GLU:OE2	3:H:237:ARG:NH1	2.43	0.49
4:I:78:ILE:HD12	4:I:84:MET:SD	2.54	0.48
5:D:271:MET:HA	5:D:271:MET:HE2	1.95	0.48
2:E:188:ARG:HH12	2:E:458:ASP:CG	2.21	0.48
3:H:6:LEU:HA	3:H:33:PHE:CZ	2.49	0.48
1:A:376:ASN:O	1:A:393:GLN:HA	2.14	0.48
5:D:81:MET:HE3	5:D:272:PHE:HB2	1.97	0.47
2:E:199:PRO:HB3	4:I:2:GLU:HB3	1.97	0.47
1:A:55:ARG:NH1	1:A:198:GLU:OE2	2.40	0.47
5:C:21:THR:CG2	5:C:27:GLY:HA2	2.44	0.47
2:F:362:ASP:OD2	2:F:385:TYR:OH	2.29	0.47
3:H:249:LYS:HE3	3:H:250:PHE:CZ	2.49	0.47
3:H:253:LYS:HB2	3:H:272:TRP:CD1	2.49	0.47
5:D:18:PHE:HE2	5:D:68:MET:SD	2.38	0.47
3:H:126:LEU:HD13	3:H:186:ILE:HD12	1.96	0.46
5:C:62:ARG:NH1	5:C:348:ASP:OD1	2.48	0.46
1:A:280:GLU:HG3	1:A:285:ASN:O	2.15	0.46
5:C:166:MET:CE	5:C:217:MET:HE1	2.45	0.46
1:B:186:LYS:HE3	5:C:297:PHE:CG	2.51	0.46
4:I:231:TYR:CE2	4:I:235:LYS:HE3	2.51	0.45
2:F:269:LEU:HD22	2:F:273:MET:HE2	1.98	0.45
3:H:24:VAL:HG22	3:H:104:ILE:HG21	1.97	0.45
2:F:512:VAL:N	2:F:513:PRO:HD2	2.31	0.45
1:B:260:PHE:CZ	1:B:272:PHE:HB2	2.51	0.45
4:J:391:LEU:C	4:J:391:LEU:HD13	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:237:ARG:HD2	3:G:304:TYR:CG	2.51	0.45
3:G:222:GLN:O	3:G:223:ASN:C	2.60	0.45
3:G:147:ARG:NH1	1:B:354:SER:O	2.50	0.45
1:A:192:HIS:ND1	1:A:194:THR:HB	2.31	0.44
1:A:409:HIS:O	1:A:410:GLN:C	2.60	0.44
1:B:70:LEU:HD22	1:B:156:ILE:CD1	2.47	0.44
3:H:100:PHE:CE2	3:H:104:ILE:HD11	2.53	0.44
1:B:132:GLY:HA2	1:B:260:PHE:CD1	2.53	0.44
1:B:206:ASP:OD1	1:B:208:THR:OG1	2.34	0.44
3:H:206:ILE:HG21	3:H:228:VAL:HG11	1.99	0.44
5:C:188:HIS:CE1	5:C:215:THR:OG1	2.70	0.44
1:A:354:SER:O	3:H:147:ARG:NH1	2.51	0.44
5:C:286:HIS:CD2	5:C:286:HIS:H	2.35	0.44
1:A:66:VAL:HB	1:A:76:ILE:HB	2.00	0.44
5:C:21:THR:HG22	5:C:27:GLY:HA2	1.99	0.43
1:A:382:GLY:O	1:A:399:GLY:HA2	2.18	0.43
2:E:326:LYS:HA	2:E:326:LYS:HD2	1.79	0.43
1:B:455:ASP:O	1:B:456:ALA:C	2.60	0.43
5:D:34:MET:HE3	5:D:75:GLU:CD	2.43	0.43
4:I:235:LYS:O	4:I:264:TYR:OH	2.31	0.43
2:E:435:LEU:N	2:E:435:LEU:HD12	2.34	0.43
3:G:147:ARG:HG2	3:G:147:ARG:HH21	1.84	0.42
2:F:366:TRP:CZ3	2:F:417:ARG:HD2	2.54	0.42
2:E:196:MET:HE1	2:E:208:VAL:HG22	2.01	0.42
1:B:432:THR:HG22	1:B:433:SER:H	1.84	0.42
2:E:408:LEU:HD12	2:E:512:VAL:HG11	2.01	0.42
1:A:432:THR:HG22	1:A:433:SER:H	1.84	0.42
2:F:403:LEU:HD13	2:F:420:THR:HG23	2.01	0.42
2:F:435:LEU:HD12	2:F:435:LEU:N	2.34	0.42
5:C:271:MET:HE2	5:C:271:MET:HA	2.01	0.42
1:A:444:PRO:O	1:A:447:SER:OG	2.30	0.42
1:A:318:VAL:HG12	1:A:318:VAL:O	2.20	0.42
5:C:125:HIS:CD2	5:C:125:HIS:C	2.97	0.42
2:E:351:THR:O	2:E:352:GLU:HB2	2.20	0.42
3:G:147:ARG:NH2	1:B:350:LEU:HD23	2.35	0.42
3:G:186:ILE:O	3:G:187:MET:C	2.63	0.42
1:B:68:LEU:HD22	1:B:397:TRP:CH2	2.55	0.42
2:F:466:LYS:HA	2:F:471:VAL:HG22	2.02	0.42
3:H:192:LEU:N	3:H:192:LEU:HD23	2.35	0.42
1:A:211:ARG:NH1	1:A:280:GLU:CD	2.77	0.42
1:A:276:LEU:C	1:A:276:LEU:HD23	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:287:ASP:OD1	3:G:287:ASP:N	2.53	0.42
2:F:515:VAL:HG13	2:F:516:LEU:HD23	2.02	0.41
3:G:202:GLU:HA	3:G:300:LEU:HD13	2.02	0.41
3:H:267:LYS:HG2	3:H:268:GLU:N	2.34	0.41
5:D:93:TYR:CZ	5:D:122:PHE:HA	2.55	0.41
5:D:125:HIS:CD2	5:D:125:HIS:C	2.97	0.41
1:A:194:THR:HG23	1:A:305:MET:HE1	2.01	0.41
2:E:237:ILE:O	2:E:297:LYS:HE3	2.20	0.41
4:I:326:VAL:N	4:I:327:PRO:CD	2.82	0.41
1:B:56:PHE:CD1	1:B:65:ARG:HD3	2.56	0.41
5:D:31:SER:OG	7:D:1000:PO4:O4	2.27	0.41
5:C:245:LEU:HG	5:C:321:ILE:HD11	2.02	0.41
1:A:394:THR:HG22	1:A:396:LEU:CD1	2.51	0.41
1:A:175:GLU:HG2	1:A:178:VAL:HG21	2.02	0.41
3:H:118:PHE:CE1	5:D:280:ASN:HB3	2.56	0.41
3:H:186:ILE:O	3:H:187:MET:C	2.63	0.41
5:D:273:LYS:NZ	7:D:1000:PO4:O4	2.48	0.41
1:B:47:LEU:C	1:B:47:LEU:HD23	2.46	0.41
4:I:5:ALA:HB2	4:I:48:PHE:CE2	2.56	0.40
1:B:211:ARG:NH1	1:B:280:GLU:CD	2.79	0.40
2:F:269:LEU:HD12	2:F:274:HIS:CE1	2.56	0.40
2:F:269:LEU:HB2	2:F:274:HIS:NE2	2.36	0.40
2:E:185:GLN:HG2	2:E:460:PRO:HD3	2.02	0.40
2:E:269:LEU:HD22	2:E:273:MET:HE2	2.04	0.40
4:I:186:ILE:HG12	4:I:237:PHE:CG	2.56	0.40
4:J:326:VAL:N	4:J:327:PRO:CD	2.84	0.40
4:I:229:ILE:HB	4:I:230:PRO:HD3	2.03	0.40
1:B:93:VAL:HG11	1:B:106:LEU:HD11	2.04	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	433/721 (60%)	414 (96%)	19 (4%)	0	100	100
1	B	433/721 (60%)	412 (95%)	21 (5%)	0	100	100
2	E	343/523 (66%)	334 (97%)	7 (2%)	2 (1%)	21	38
2	F	343/523 (66%)	339 (99%)	4 (1%)	0	100	100
3	G	303/322 (94%)	292 (96%)	11 (4%)	0	100	100
3	H	303/322 (94%)	294 (97%)	9 (3%)	0	100	100
4	I	385/452 (85%)	361 (94%)	22 (6%)	2 (0%)	24	43
4	J	385/452 (85%)	366 (95%)	18 (5%)	1 (0%)	36	55
5	C	318/368 (86%)	311 (98%)	7 (2%)	0	100	100
5	D	318/368 (86%)	314 (99%)	4 (1%)	0	100	100
All	All	3564/4772 (75%)	3437 (96%)	122 (3%)	5 (0%)	49	68

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	190	ASN
4	I	313	VAL
4	I	393	MET
2	E	184	PRO
4	J	151	ALA

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	384/626 (61%)	376 (98%)	8 (2%)	47	74
1	B	384/626 (61%)	378 (98%)	6 (2%)	55	79
2	E	302/444 (68%)	294 (97%)	8 (3%)	40	68
2	F	302/444 (68%)	294 (97%)	8 (3%)	40	68
3	G	260/274 (95%)	256 (98%)	4 (2%)	57	80
3	H	260/274 (95%)	252 (97%)	8 (3%)	35	62

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
4	I	345/398 (87%)	343 (99%)	2 (1%)	78 91
4	J	345/398 (87%)	342 (99%)	3 (1%)	70 87
5	C	274/312 (88%)	273 (100%)	1 (0%)	84 93
5	D	274/312 (88%)	267 (97%)	7 (3%)	40 68
All	All	3130/4108 (76%)	3075 (98%)	55 (2%)	51 77

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	PHE
1	A	65	ARG
1	A	117	LEU
1	A	151	VAL
1	A	161	ILE
1	A	180	VAL
1	A	223	ARG
1	A	422	ARG
2	E	185	GLN
2	E	241	THR
2	E	246	GLU
2	E	328	SER
2	E	386	LEU
2	E	415	MET
2	E	487	LEU
2	E	518	VAL
3	G	24	VAL
3	G	63	VAL
3	G	104	ILE
3	G	160	SER
4	I	18	ASP
4	I	118	VAL
1	B	155	VAL
1	B	175	GLU
1	B	196	CYS
1	B	199	ASP
1	B	422	ARG
1	B	427	PRO
2	F	246	GLU
2	F	386	LEU
2	F	396	PRO
2	F	415	MET

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Mol	Chain	Res	Type
2	F	448	GLN
2	F	487	LEU
2	F	495	GLU
2	F	518	VAL
3	H	8	GLU
3	H	11	LYS
3	H	13	GLN
3	H	83	TYR
3	H	87	SER
3	H	142	VAL
3	H	192	LEU
3	H	300	LEU
4	J	128	LEU
4	J	345	GLN
4	J	404	ILE
5	C	350	VAL
5	D	91	GLU
5	D	161	SER
5	D	165	ILE
5	D	202	GLU
5	D	327	ASN
5	D	328	ILE
5	D	333	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	216	GLN
1	A	337	HIS
1	A	340	HIS
2	E	185	GLN
2	E	265	GLN
3	G	244	GLN
4	I	405	GLN
1	B	160	ASN
1	B	216	GLN
2	F	455	ASN
3	H	208	ASN
4	J	363	GLN
4	J	403	ASN
5	C	132	ASN
5	C	277	GLN

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Mol	Chain	Res	Type
5	D	72	GLN
5	D	261	HIS
5	D	277	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)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\#\|Z\| > 2$
7	PO4	D	1000	-	4,4,4	1.33	1 (25%)	6,6,6	0.75	0
7	PO4	C	1000	-	4,4,4	1.64	1 (25%)	6,6,6	0.56	0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	1000	PO4	P-O1	3.06	1.57	1.50
7	D	1000	PO4	P-O1	2.43	1.56	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	1000	PO4	2	0

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.

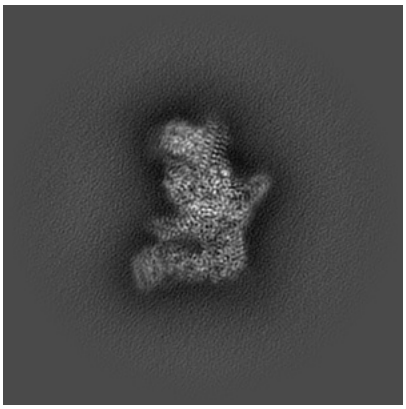
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-75149. These allow visual inspection of the internal detail of the map and identification of artifacts.

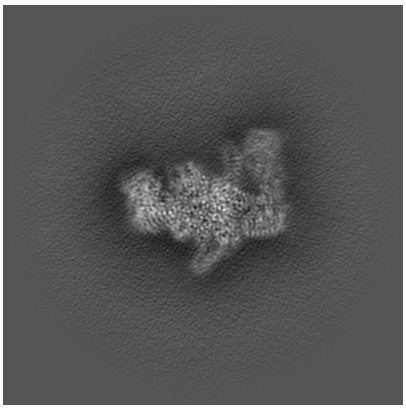
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

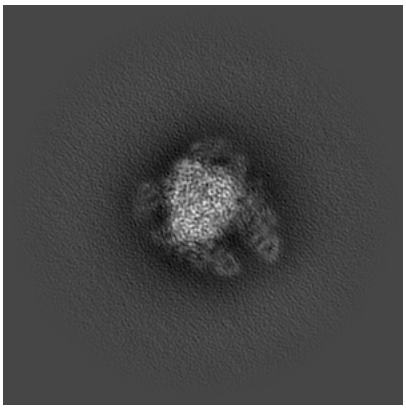
6.1.1 Primary map



X

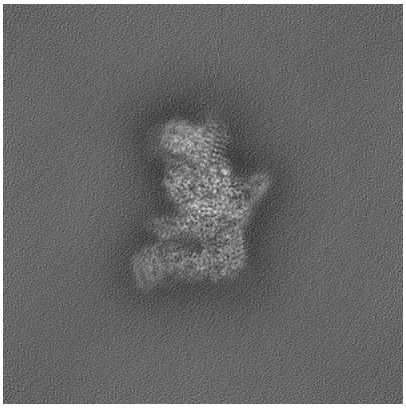


Y

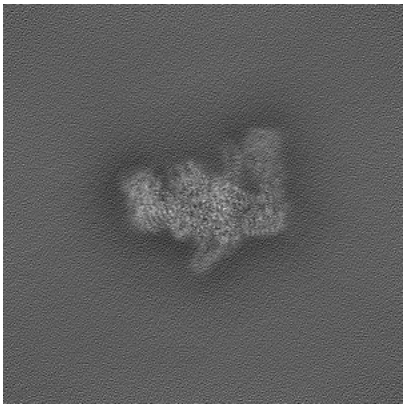


Z

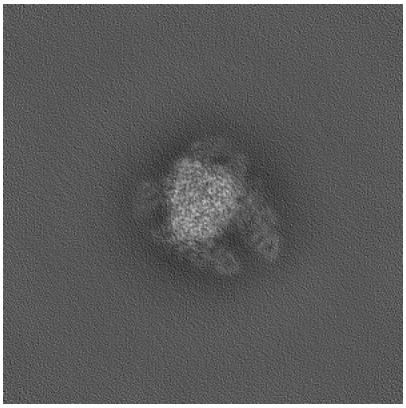
6.1.2 Raw map



X



Y

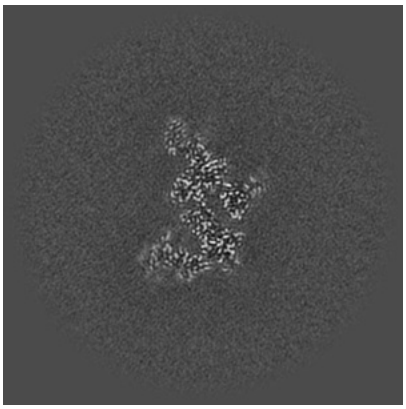


Z

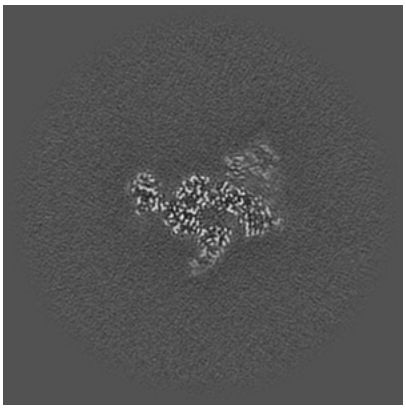
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

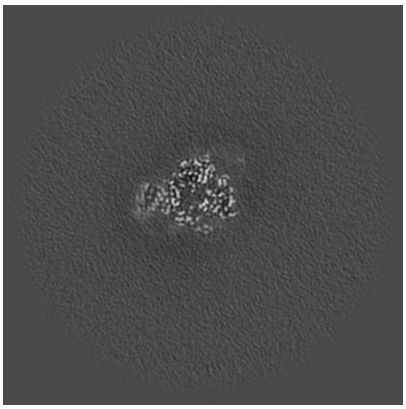
6.2.1 Primary map



X Index: 200

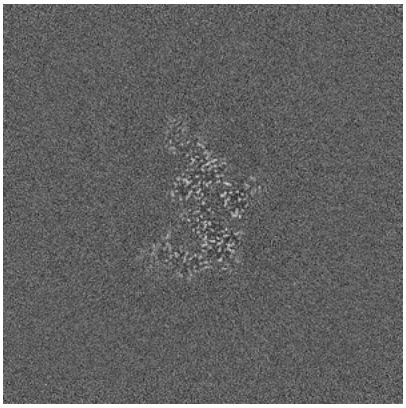


Y Index: 200

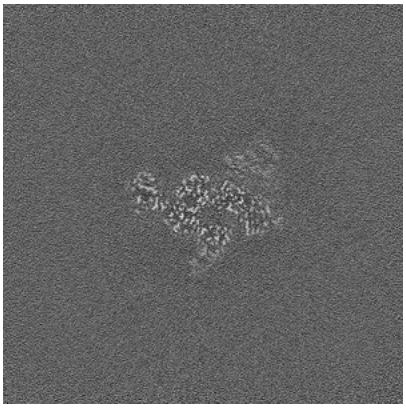


Z Index: 200

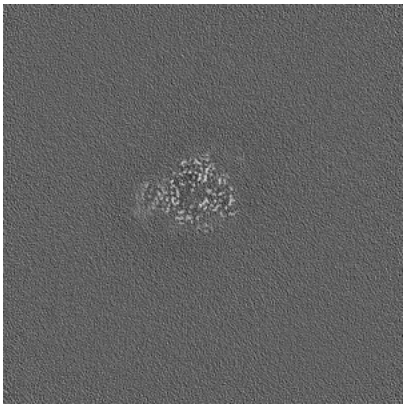
6.2.2 Raw map



X Index: 200



Y Index: 200

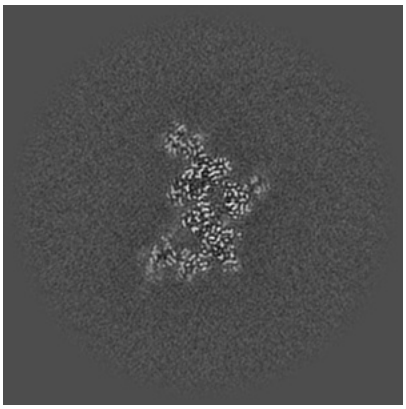


Z Index: 200

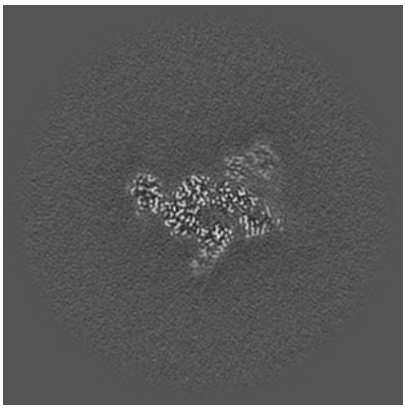
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

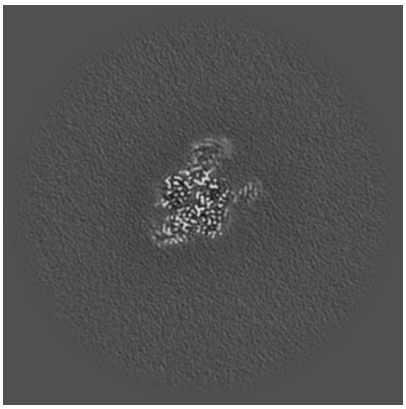
6.3.1 Primary map



X Index: 202

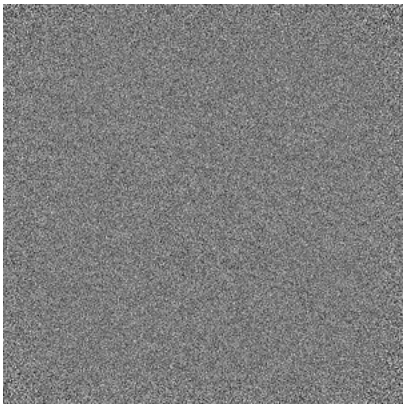


Y Index: 201

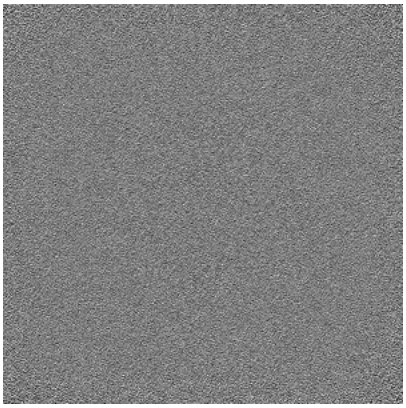


Z Index: 220

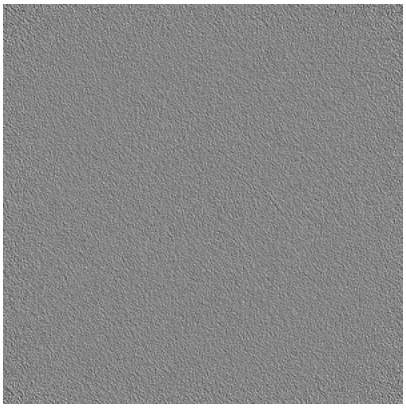
6.3.2 Raw map



X Index: 0



Y Index: 0

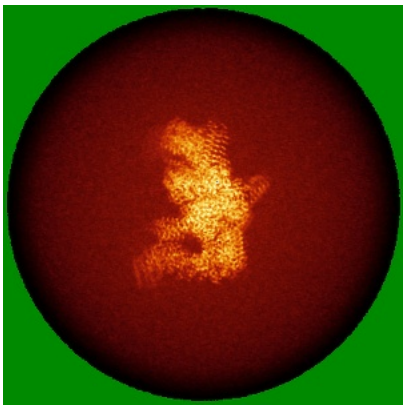


Z Index: 0

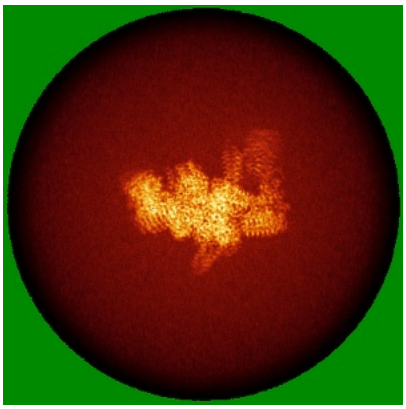
The images above show the largest variance slices of the map in three orthogonal directions.

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

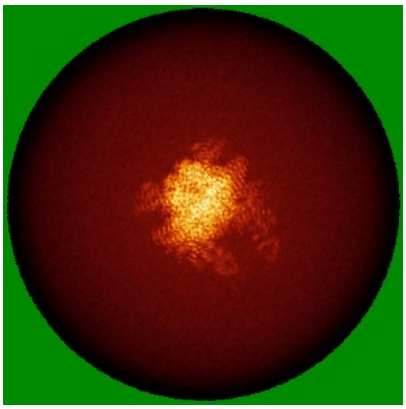
6.4.1 Primary map



X

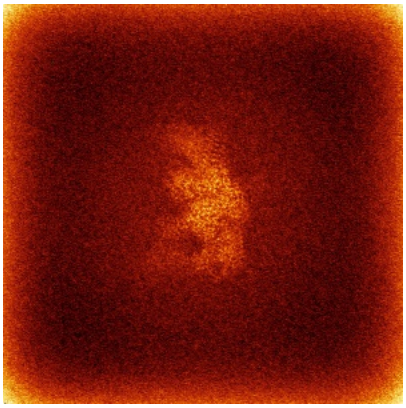


Y

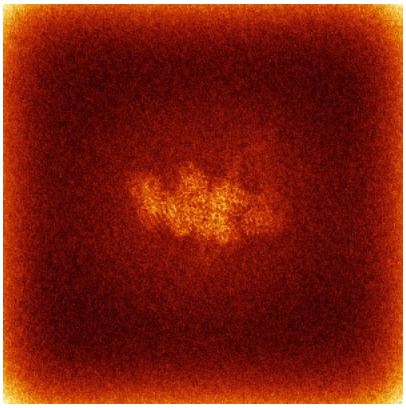


Z

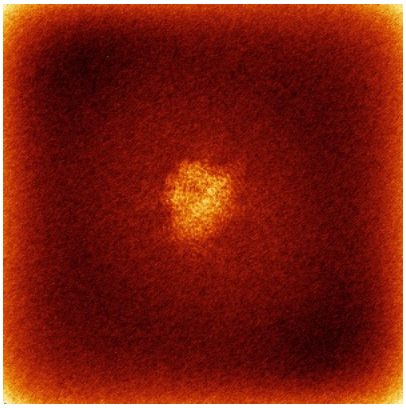
6.4.2 Raw map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views (i)

6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.165. 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.

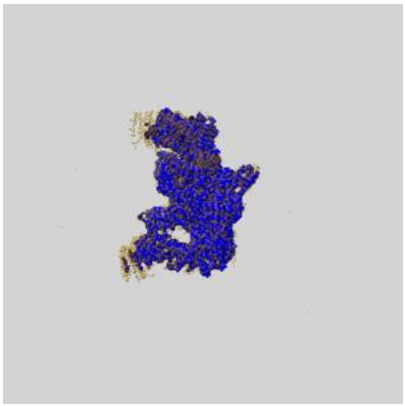
6.6 Mask visualisation (i)

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

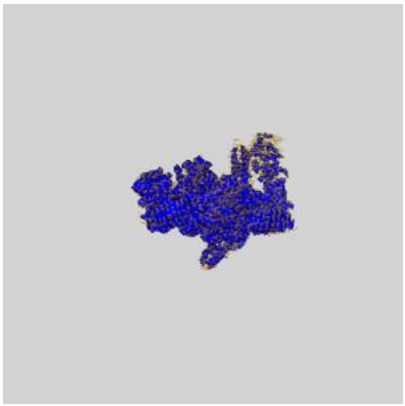
A mask typically either:

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

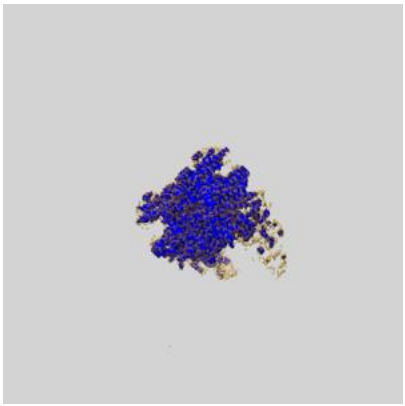
6.6.1 emd_75149_msk_1.map (i)



X



Y

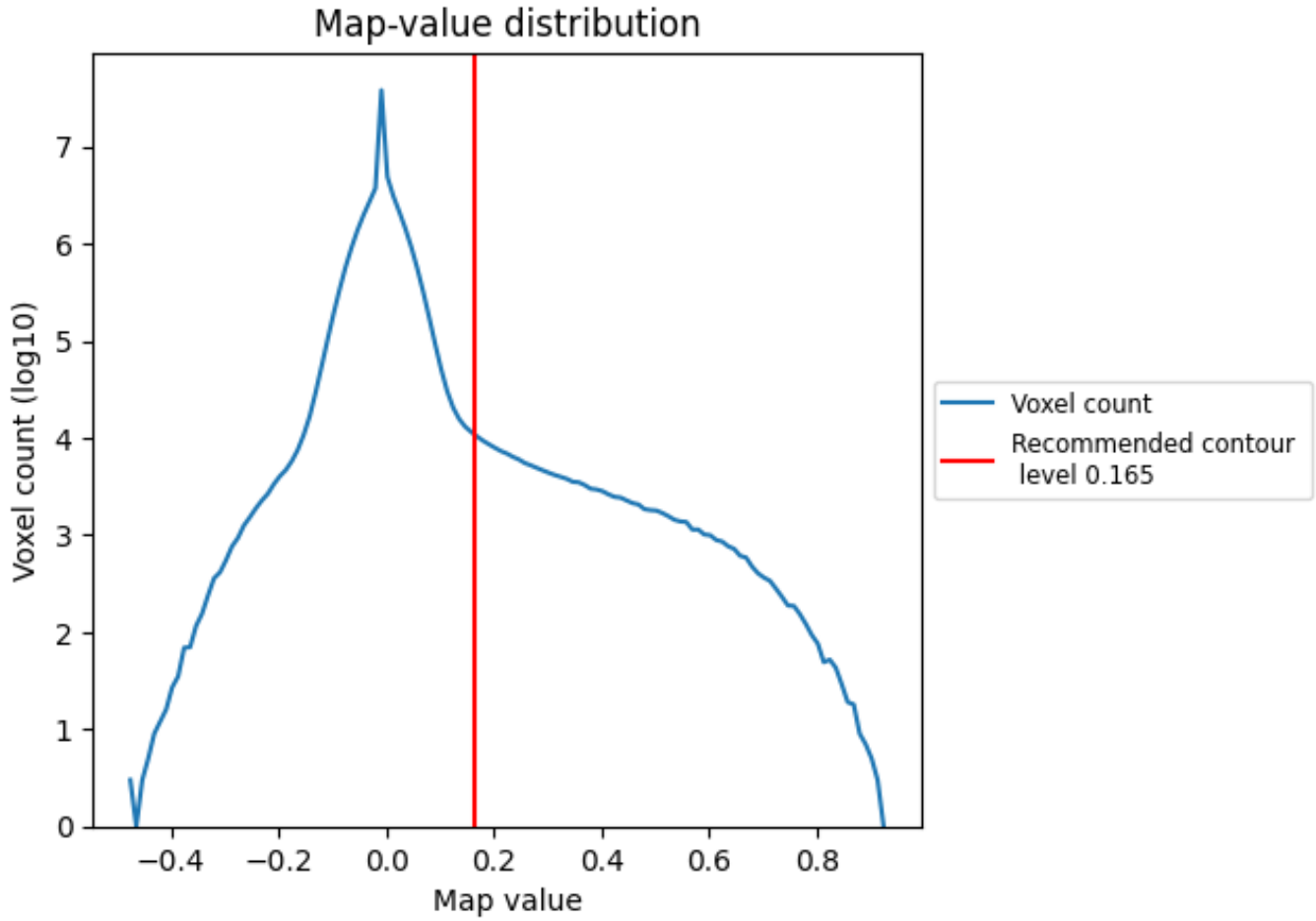


Z

7 Map analysis

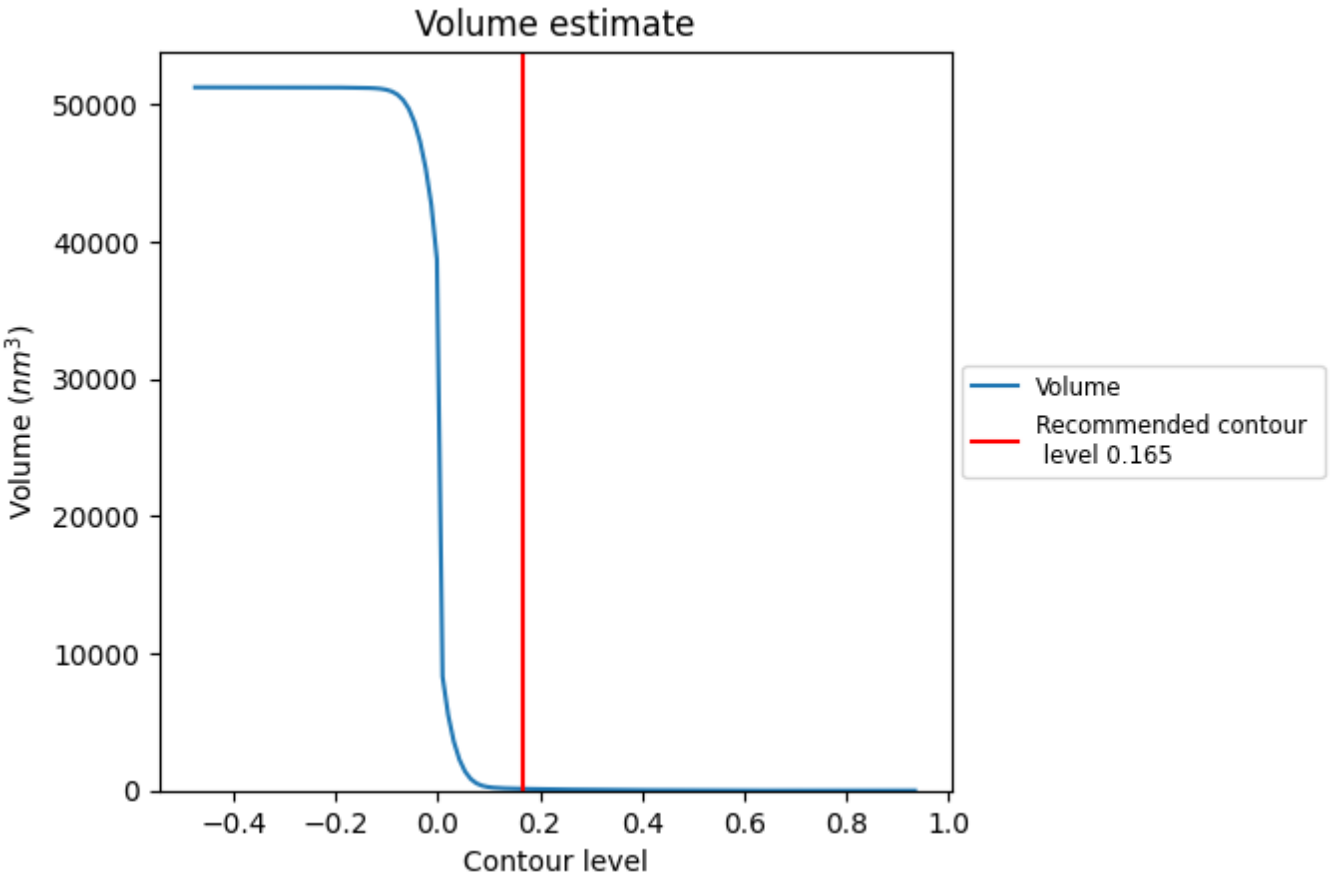
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

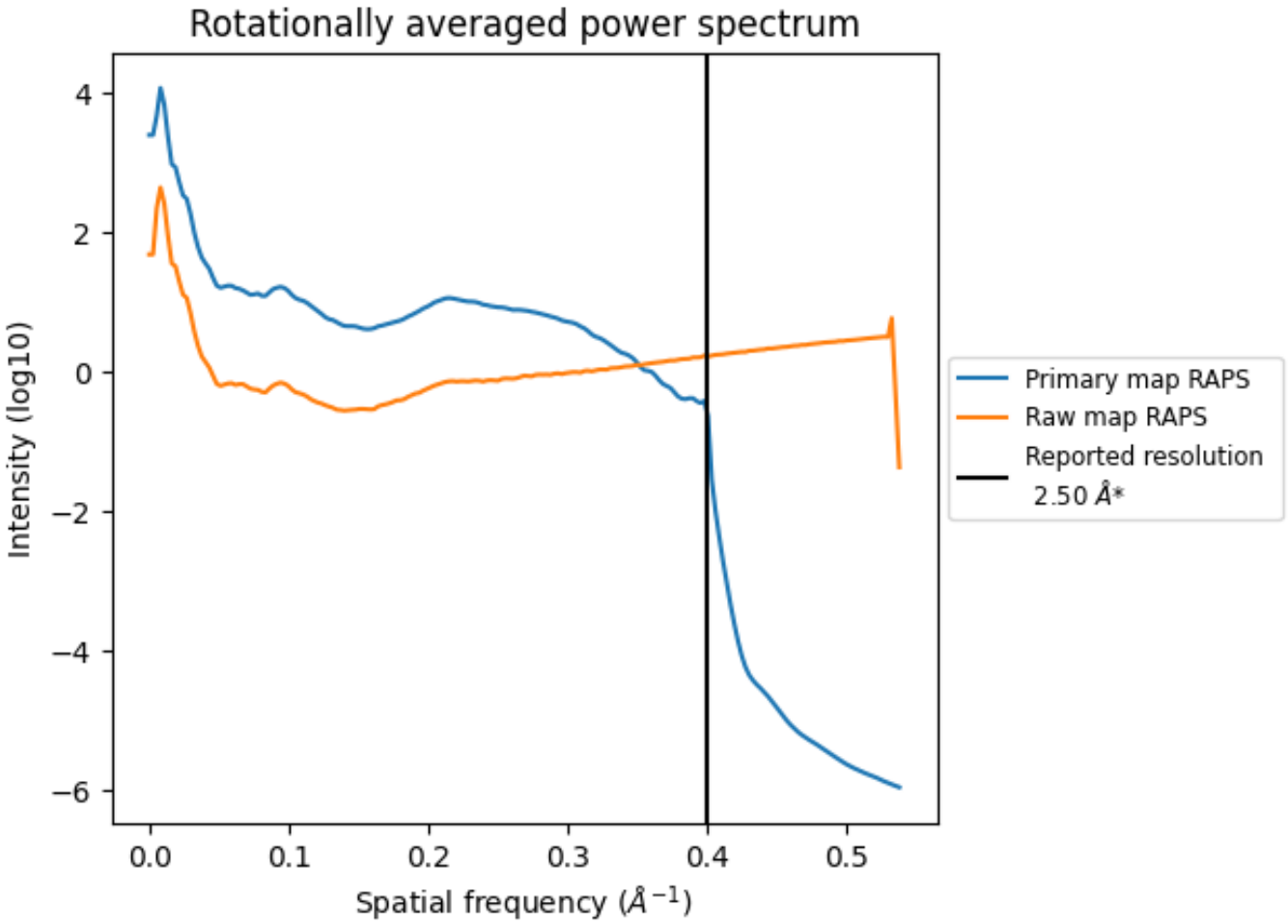
7.2 Volume estimate



The volume at the recommended contour level is 129 nm^3; this corresponds to an approximate mass of 116 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)

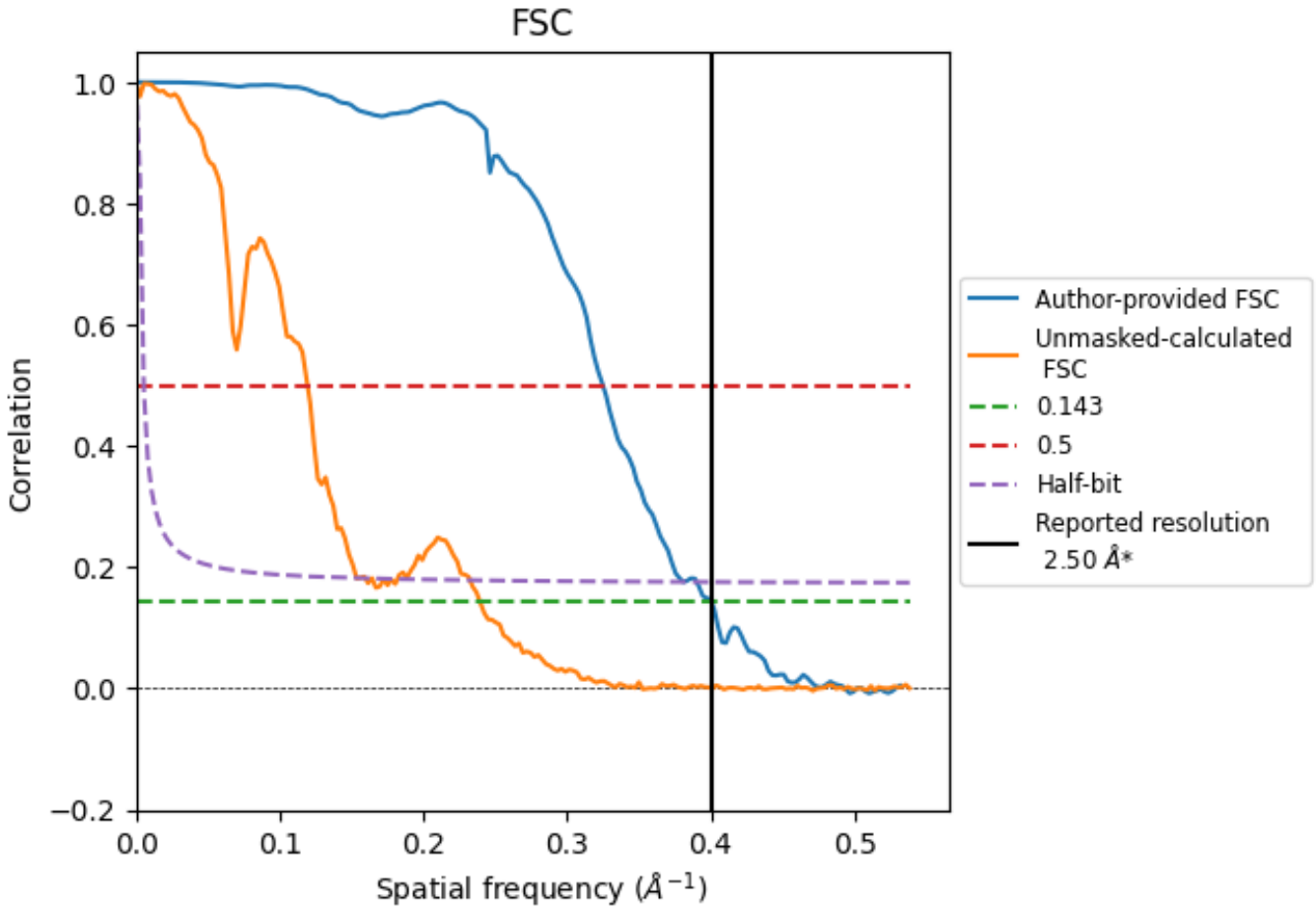


*Reported resolution corresponds to spatial frequency of 0.400 $Å^{-1}$

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.400 Å^{-1}

8.2 Resolution estimates (i)

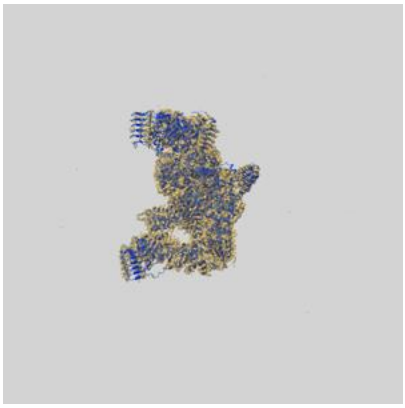
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.50	-	-
Author-provided FSC curve	2.50	3.08	2.63
Unmasked-calculated*	4.18	8.39	6.52

*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 4.18 differs from the reported value 2.5 by more than 10 %

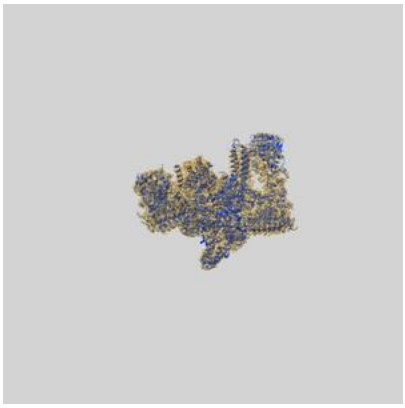
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-75149 and PDB model 10GF. Per-residue inclusion information can be found in section 3 on page 8.

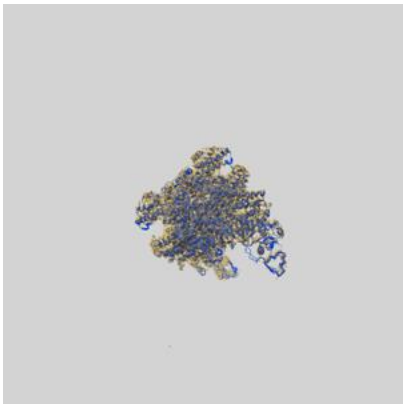
9.1 Map-model overlay



X



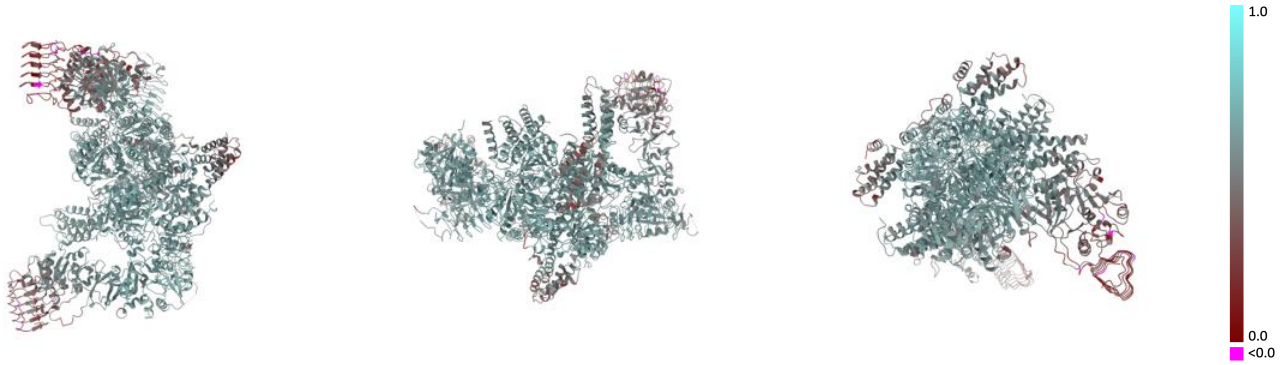
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.165 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.

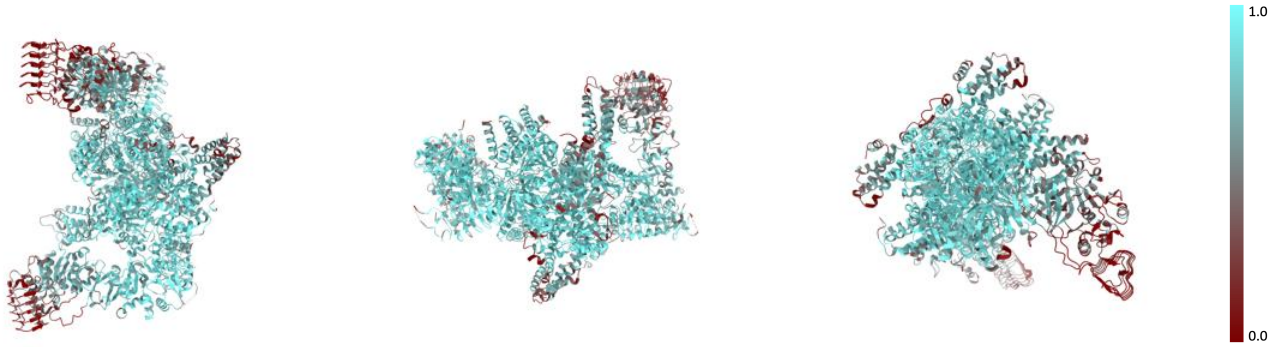
9.2 Q-score mapped to coordinate model (i)



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.

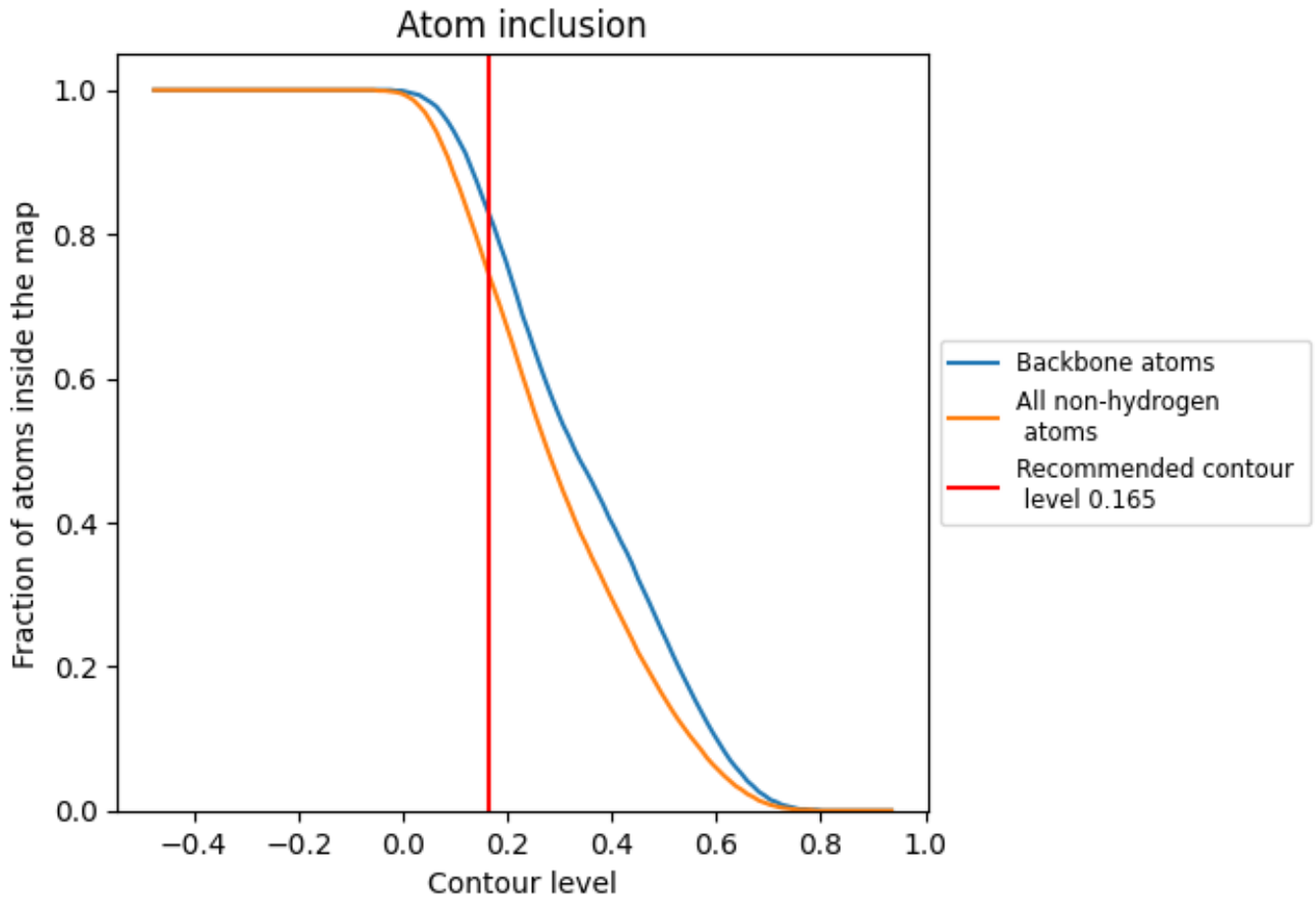
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.165).

9.4 Atom inclusion (i)



At the recommended contour level, 83% of all backbone atoms, 74% 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 (0.165) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7430	0.5540
A	0.8770	0.6090
B	0.7930	0.5760
C	0.8760	0.6060
D	0.8900	0.6120
E	0.8620	0.6080
F	0.7820	0.5870
G	0.7580	0.5620
H	0.7590	0.5540
I	0.5450	0.4700
J	0.3400	0.3750

