



Full wwPDB EM Validation Report ⓘ

Sep 27, 2023 – 05:40 PM EDT

PDB ID : 8HCX
EMDB ID : EMD-34667
Title : Cryo-EM structure of Endothelin1-bound ETBR-Gq complex
Authors : Yuan, Q.; Jiang, Y.; Xu, H.E.; Ji, Y.; Duan, J.
Deposited on : 2022-11-03
Resolution : 3.50 Å (reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev50
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

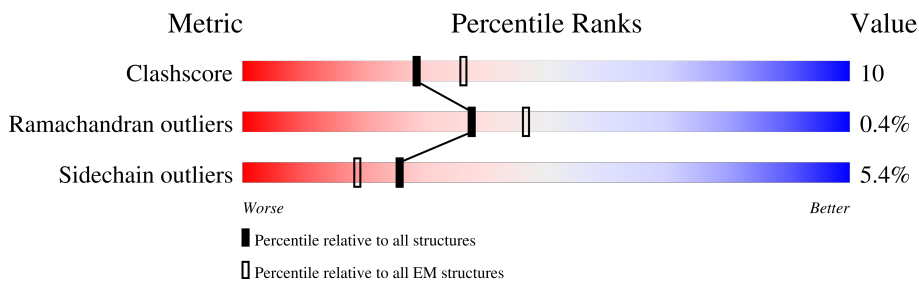
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	246	
2	B	377	
3	C	603	
4	D	21	
5	E	285	
6	G	71	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8933 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 Guanine nucleotide-binding protein G(q) subunit alpha-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	219	1780	1123	311	339	7	0	0

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	338	2596	1604	467	504	21	0	0

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	MET	-	initiating methionine	UNP P62873
B	-4	HIS	-	expression tag	UNP P62873
B	-3	HIS	-	expression tag	UNP P62873
B	-2	HIS	-	expression tag	UNP P62873
B	-1	HIS	-	expression tag	UNP P62873
B	0	HIS	-	expression tag	UNP P62873
B	1	HIS	-	expression tag	UNP P62873
B	2	GLY	-	expression tag	UNP P62873
B	3	SER	-	expression tag	UNP P62873
B	4	LEU	-	expression tag	UNP P62873
B	5	LEU	-	expression tag	UNP P62873
B	6	GLN	-	expression tag	UNP P62873
B	346	GLY	-	expression tag	UNP P62873
B	347	SER	-	expression tag	UNP P62873
B	348	SER	-	expression tag	UNP P62873
B	349	GLY	-	expression tag	UNP P62873
B	350	GLY	-	expression tag	UNP P62873
B	351	GLY	-	expression tag	UNP P62873
B	352	GLY	-	expression tag	UNP P62873
B	353	SER	-	expression tag	UNP P62873

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Chain	Residue	Modelled	Actual	Comment	Reference
B	354	GLY	-	expression tag	UNP P62873
B	355	GLY	-	expression tag	UNP P62873
B	356	GLY	-	expression tag	UNP P62873
B	357	GLY	-	expression tag	UNP P62873
B	358	SER	-	expression tag	UNP P62873
B	359	SER	-	expression tag	UNP P62873
B	360	GLY	-	expression tag	UNP P62873
B	361	VAL	-	expression tag	UNP P62873
B	362	SER	-	expression tag	UNP P62873
B	363	GLY	-	expression tag	UNP P62873
B	364	TRP	-	expression tag	UNP P62873
B	365	ARG	-	expression tag	UNP P62873
B	366	LEU	-	expression tag	UNP P62873
B	367	PHE	-	expression tag	UNP P62873
B	368	LYS	-	expression tag	UNP P62873
B	369	LYS	-	expression tag	UNP P62873
B	370	ILE	-	expression tag	UNP P62873
B	371	SER	-	expression tag	UNP P62873

- Molecule 3 is a protein called Endothelin receptor type B,Oplophorus-luciferin 2-monooxygenase catalytic subunit chimera.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	292	2190	1467	334	370	19	0	0

There are 77 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-20	MET	-	initiating methionine	UNP P24530
C	-19	ASP	-	expression tag	UNP P24530
C	-18	SER	-	expression tag	UNP P24530
C	-17	LYS	-	expression tag	UNP P24530
C	-16	GLY	-	expression tag	UNP P24530
C	-15	SER	-	expression tag	UNP P24530
C	-14	SER	-	expression tag	UNP P24530
C	-13	GLN	-	expression tag	UNP P24530
C	-12	LYS	-	expression tag	UNP P24530
C	-11	GLY	-	expression tag	UNP P24530
C	-10	SER	-	expression tag	UNP P24530
C	-9	ARG	-	expression tag	UNP P24530
C	-8	LEU	-	expression tag	UNP P24530

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-7	LEU	-	expression tag	UNP P24530
C	-6	LEU	-	expression tag	UNP P24530
C	-5	LEU	-	expression tag	UNP P24530
C	-4	LEU	-	expression tag	UNP P24530
C	-3	VAL	-	expression tag	UNP P24530
C	-2	VAL	-	expression tag	UNP P24530
C	-1	SER	-	expression tag	UNP P24530
C	0	ASN	-	expression tag	UNP P24530
C	1	LEU	-	expression tag	UNP P24530
C	2	LEU	-	expression tag	UNP P24530
C	3	LEU	-	expression tag	UNP P24530
C	4	CYS	-	expression tag	UNP P24530
C	5	GLN	-	expression tag	UNP P24530
C	6	GLY	-	expression tag	UNP P24530
C	7	VAL	-	expression tag	UNP P24530
C	8	VAL	-	expression tag	UNP P24530
C	9	SER	-	expression tag	UNP P24530
C	10	ASP	-	expression tag	UNP P24530
C	11	TYR	-	expression tag	UNP P24530
C	12	LYS	-	expression tag	UNP P24530
C	13	ASP	-	expression tag	UNP P24530
C	14	ASP	-	expression tag	UNP P24530
C	15	ASP	-	expression tag	UNP P24530
C	16	ASP	-	expression tag	UNP P24530
C	17	VAL	-	expression tag	UNP P24530
C	18	ASP	-	expression tag	UNP P24530
C	19	HIS	-	expression tag	UNP P24530
C	20	HIS	-	expression tag	UNP P24530
C	21	HIS	-	expression tag	UNP P24530
C	22	HIS	-	expression tag	UNP P24530
C	23	HIS	-	expression tag	UNP P24530
C	24	HIS	-	expression tag	UNP P24530
C	25	HIS	-	expression tag	UNP P24530
C	26	HIS	-	expression tag	UNP P24530
C	425	VAL	THR	conflict	UNP Q9GV45
C	429	GLU	ALA	conflict	UNP Q9GV45
C	436	GLU	GLN	conflict	UNP Q9GV45
C	440	ALA	GLY	conflict	UNP Q9GV45
C	443	LEU	GLN	conflict	UNP Q9GV45
C	452	VAL	LEU	conflict	UNP Q9GV45
C	456	LEU	PHE	conflict	UNP Q9GV45
C	458	ASN	ALA	conflict	UNP Q9GV45

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Chain	Residue	Modelled	Actual	Comment	Reference
C	460	ALA	GLY	conflict	UNP Q9GV45
C	468	ARG	LYS	conflict	UNP Q9GV45
C	469	ILE	VAL	conflict	UNP Q9GV45
C	471	ARG	LEU	conflict	UNP Q9GV45
C	476	ALA	GLY	conflict	UNP Q9GV45
C	479	ILE	ALA	conflict	UNP Q9GV45
C	492	ALA	GLY	conflict	UNP Q9GV45
C	493	ASP	PHE	conflict	UNP Q9GV45
C	496	ALA	GLY	conflict	UNP Q9GV45
C	497	GLN	LEU	conflict	UNP Q9GV45
C	500	GLU	MET	conflict	UNP Q9GV45
C	501	VAL	ILE	conflict	UNP Q9GV45
C	515	VAL	ILE	conflict	UNP Q9GV45
C	518	PRO	HIS	conflict	UNP Q9GV45
C	532	LEU	ILE	conflict	UNP Q9GV45
C	533	ASN	ASP	conflict	UNP Q9GV45
C	540	GLU	PRO	conflict	UNP Q9GV45
C	549	LYS	GLN	conflict	UNP Q9GV45
C	563	ILE	TYR	conflict	UNP Q9GV45
C	569	THR	ASN	conflict	UNP Q9GV45
C	574	MET	LEU	conflict	UNP Q9GV45
C	582	SER	GLY	conflict	UNP Q9GV45

- Molecule 4 is a protein called Endothelin-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	21	164	104	24	32	4	0	0

- Molecule 5 is a protein called scFv16.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	E	232	1770	1126	292	342	10	0	0

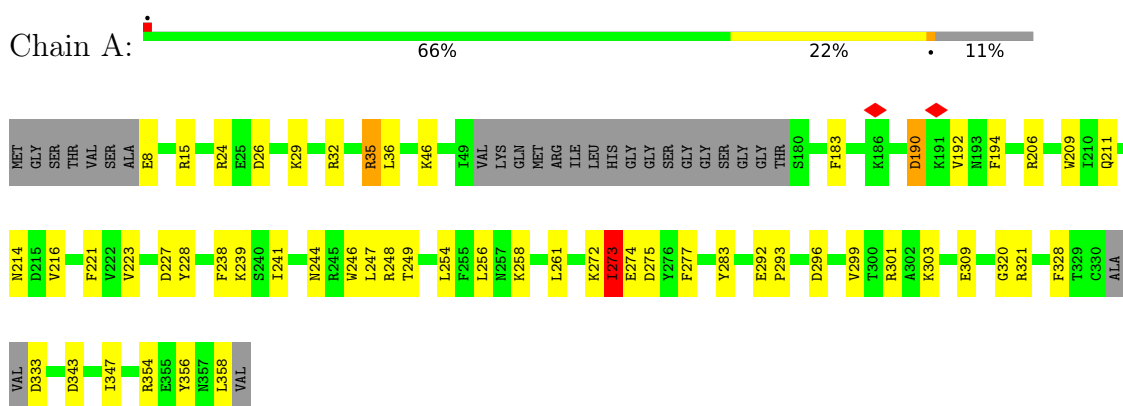
- Molecule 6 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	G	57	433	271	77	83	2	0	0

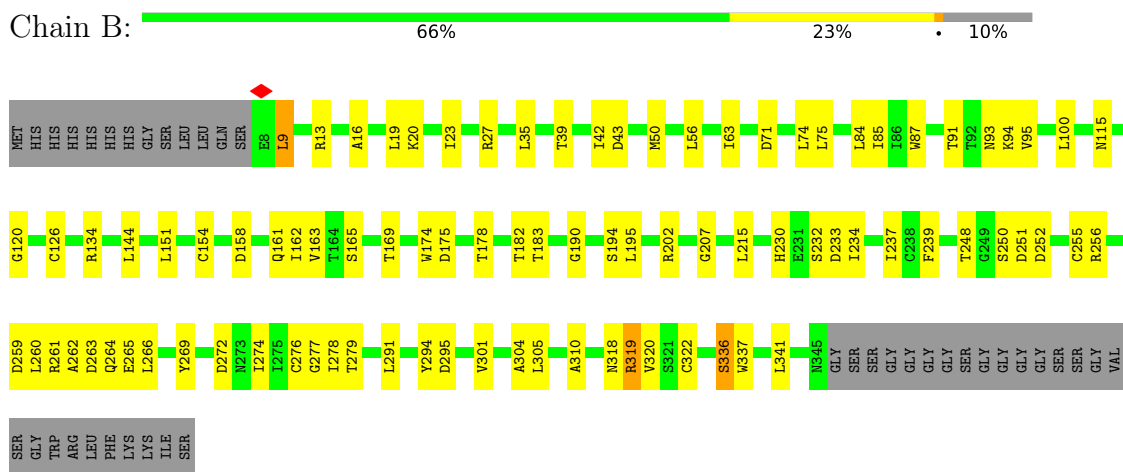
3 Residue-property plots

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.

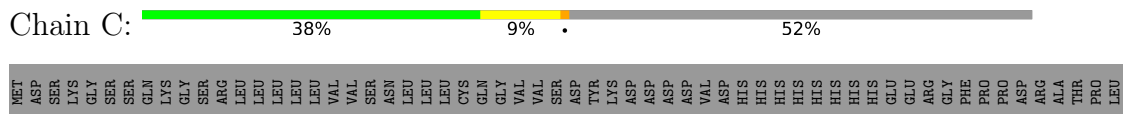
- Molecule 1: Guanine nucleotide-binding protein G(q) subunit alpha-1



- Molecule 2: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1



- Molecule 3: Endothelin receptor type B, Ophiophorus-luciferin 2-monooxygenase catalytic subunit chimera



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	373655	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	1.880	Depositor
Minimum map value	-0.003	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.026	Depositor
Recommended contour level	0.013	Depositor
Map size (\AA)	266.24, 266.24, 266.24	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.04, 1.04, 1.04	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	0.26	0/1812	0.55	1/2440 (0.0%)
2	B	0.25	0/2643	0.54	0/3583
3	C	0.26	0/2242	0.46	1/3074 (0.0%)
4	D	0.23	0/168	0.38	0/227
5	E	0.26	0/1814	0.53	0/2461
6	G	0.24	0/439	0.45	0/594
All	All	0.25	0/9118	0.51	2/12379 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	ASP	CB-CG-OD1	7.06	124.65	118.30
3	C	149	LEU	CA-CB-CG	5.53	128.01	115.30

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	1780	0	1725	32	0
2	B	2596	0	2505	63	0
3	C	2190	0	2172	46	0
4	D	164	0	141	6	0
5	E	1770	0	1709	39	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	G	433	0	441	6	0
All	All	8933	0	8693	184	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 10.

All (184) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:94:ILE:HG13	3:C:96:ILE:H	1.52	0.73
2:B:234:ILE:HA	2:B:250:SER:HA	1.71	0.72
2:B:9:LEU:HD12	2:B:13:ARG:HH22	1.55	0.72
2:B:318:ASN:OD1	2:B:319:ARG:N	2.22	0.72
5:E:231:GLN:HE21	5:E:238:THR:HG22	1.55	0.70
2:B:84:LEU:HB2	2:B:100:LEU:HD21	1.73	0.70
2:B:85:ILE:HD11	2:B:94:LYS:HG2	1.74	0.69
3:C:365:LEU:HD21	4:D:18:ASP:HA	1.75	0.69
2:B:91:THR:OG1	2:B:93:ASN:OD1	2.12	0.68
5:E:138:ILE:HD12	5:E:163:LYS:HG2	1.78	0.65
2:B:255:CYS:HB2	2:B:269:TYR:HB2	1.81	0.63
5:E:89:ASP:O	5:E:93:TYR:OH	2.17	0.60
3:C:181:GLN:HE22	3:C:182:LYS:HD2	1.65	0.60
3:C:94:ILE:HG13	3:C:96:ILE:N	2.17	0.60
5:E:38:GLN:HB3	5:E:92:MET:HB3	1.83	0.59
2:B:163:VAL:HG12	2:B:195:LEU:HD11	1.82	0.59
2:B:207:GLY:HA3	2:B:234:ILE:HD11	1.84	0.59
2:B:232:SER:HB3	2:B:252:ASP:HB3	1.84	0.59
5:E:142:GLN:NE2	5:E:241:ALA:O	2.35	0.59
3:C:180:ILE:O	3:C:184:SER:N	2.35	0.59
5:E:180:ARG:NH1	5:E:222:GLU:O	2.37	0.58
2:B:27:ARG:O	2:B:264:GLN:NE2	2.37	0.57
2:B:230:HIS:CE1	2:B:256:ARG:HG3	2.39	0.57
3:C:254:ILE:HD13	4:D:12:VAL:HB	1.85	0.57
1:A:292:GLU:HG2	1:A:293:PRO:HD2	1.86	0.57
1:A:320:GLY:O	1:A:354:ARG:NH1	2.38	0.57
5:E:178:LEU:HB2	5:E:188:LEU:HD11	1.87	0.56
3:C:104:ASN:ND2	3:C:373:ASN:OD1	2.37	0.56
2:B:232:SER:OG	2:B:233:ASP:N	2.38	0.56
5:E:75:LYS:HD2	5:E:79:PHE:HZ	1.71	0.55
2:B:19:LEU:O	2:B:23:ILE:HG23	2.07	0.55
2:B:16:ALA:HB2	6:G:16:VAL:HG12	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:16:ALA:O	2:B:20:LYS:HG2	2.07	0.54
2:B:336:SER:OG	2:B:337:TRP:N	2.39	0.54
2:B:85:ILE:HG12	2:B:87:TRP:NE1	2.22	0.53
3:C:181:GLN:HA	3:C:184:SER:HB3	1.89	0.53
2:B:120:GLY:HA3	2:B:151:LEU:HD23	1.90	0.53
3:C:181:GLN:NE2	3:C:182:LYS:HD2	2.24	0.53
3:C:94:ILE:HG22	3:C:362:SER:HA	1.90	0.53
3:C:100:PHE:HE2	3:C:366:VAL:HG22	1.74	0.53
2:B:301:VAL:HG12	2:B:310:ALA:HB3	1.91	0.52
3:C:332:PHE:HD1	3:C:378:ASN:HD21	1.57	0.52
3:C:274:ASP:HB3	3:C:347:LEU:HD22	1.92	0.52
1:A:211:GLN:HG2	1:A:246:TRP:CD1	2.44	0.52
2:B:262:ALA:HB1	6:G:30:VAL:HG12	1.92	0.52
3:C:185:VAL:HG13	3:C:281:TYR:OH	2.10	0.51
3:C:182:LYS:HE3	3:C:277:LEU:HD21	1.91	0.51
3:C:324:THR:HG21	3:C:389:VAL:HG21	1.93	0.51
3:C:109:CYS:O	3:C:113:VAL:HG22	2.10	0.51
3:C:277:LEU:O	3:C:281:TYR:HB2	2.11	0.51
2:B:63:ILE:HD13	2:B:341:LEU:HD22	1.92	0.51
5:E:39:ALA:HB3	5:E:42:LYS:HG3	1.92	0.50
5:E:86:ARG:NH2	5:E:88:GLU:OE1	2.43	0.50
1:A:214:ASN:HD21	1:A:246:TRP:HB3	1.76	0.50
3:C:173:MET:SD	3:C:177:VAL:HG23	2.51	0.50
2:B:194:SER:HB3	2:B:237:ILE:HG22	1.94	0.50
2:B:56:LEU:HB2	2:B:341:LEU:HB2	1.94	0.50
3:C:154:ASP:OD1	3:C:155:ILE:N	2.45	0.50
3:C:282:PHE:O	3:C:286:LEU:HB2	2.12	0.49
3:C:343:ARG:HH11	3:C:343:ARG:HG3	1.78	0.49
3:C:254:ILE:HA	4:D:16:HIS:HD2	1.78	0.49
5:E:31:PHE:CD2	5:E:97:ARG:HD2	2.48	0.49
5:E:98:SER:HB3	5:E:107:PHE:H	1.77	0.48
5:E:39:ALA:HA	5:E:91:ALA:HA	1.94	0.48
1:A:272:LYS:O	1:A:274:GLU:N	2.46	0.48
5:E:31:PHE:CZ	5:E:99:ILE:HD11	2.49	0.48
5:E:36:VAL:HG23	5:E:94:TYR:HB2	1.96	0.48
6:G:13:ARG:O	6:G:17:GLU:HG2	2.14	0.48
1:A:206:ARG:HA	1:A:209:TRP:CD1	2.49	0.48
2:B:134:ARG:NH1	5:E:197:SER:OG	2.47	0.48
2:B:276:CYS:HB2	2:B:295:ASP:HB2	1.95	0.48
2:B:85:ILE:HG23	2:B:87:TRP:HE1	1.79	0.47
1:A:272:LYS:C	1:A:274:GLU:H	2.16	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:56:LEU:HD21	2:B:75:LEU:HD13	1.96	0.47
2:B:85:ILE:HG12	2:B:87:TRP:CD1	2.50	0.47
2:B:71:ASP:N	2:B:71:ASP:OD1	2.42	0.47
5:E:72:ASP:HB3	5:E:75:LYS:HG2	1.96	0.47
1:A:296:ASP:HB3	1:A:299:VAL:HB	1.96	0.47
2:B:182:THR:HG22	2:B:183:THR:HG23	1.97	0.47
3:C:336:TRP:CZ3	3:C:339:LEU:HD23	2.50	0.47
5:E:96:VAL:HG11	5:E:107:PHE:CD2	2.50	0.47
1:A:26:ASP:OD1	1:A:29:LYS:NZ	2.48	0.46
2:B:42:ILE:HG22	2:B:43:ASP:N	2.30	0.46
3:C:126:ILE:HD11	3:C:137:ASN:HB3	1.96	0.46
1:A:283:TYR:O	1:A:301:ARG:NH2	2.35	0.46
1:A:190:ASP:O	1:A:192:VAL:HG12	2.16	0.46
3:C:149:LEU:O	3:C:153:ILE:HG22	2.15	0.46
5:E:97:ARG:NH1	5:E:108:ASP:OD1	2.48	0.46
2:B:93:ASN:OD1	2:B:93:ASN:N	2.41	0.46
3:C:277:LEU:HD23	3:C:277:LEU:HA	1.79	0.46
5:E:19:LEU:HD11	5:E:114:THR:HG21	1.97	0.46
5:E:28:PHE:HZ	5:E:78:LEU:HB2	1.81	0.46
1:A:221:PHE:CE2	1:A:241:ILE:HD11	2.51	0.46
3:C:279:SER:HA	3:C:283:CYS:HB2	1.97	0.46
2:B:255:CYS:SG	2:B:278:ILE:HG12	2.55	0.46
3:C:122:LEU:HD13	3:C:144:ALA:HB2	1.96	0.46
5:E:174:LEU:HD22	5:E:212:PHE:HD2	1.81	0.46
5:E:68:THR:HB	5:E:81:GLN:HB2	1.98	0.46
6:G:52:THR:O	6:G:52:THR:OG1	2.34	0.46
2:B:39:THR:HG21	2:B:305:LEU:HB3	1.98	0.45
2:B:126:CYS:HB3	2:B:144:LEU:HD12	1.98	0.45
1:A:356:TYR:HB2	1:A:358:LEU:HD22	1.96	0.45
2:B:202:ARG:NE	2:B:202:ARG:HA	2.31	0.45
3:C:225:ILE:HA	3:C:228:VAL:HG12	1.98	0.45
3:C:378:ASN:HA	3:C:381:ILE:HG12	1.99	0.45
2:B:291:LEU:HG	2:B:301:VAL:HG22	1.97	0.45
1:A:303:LYS:HB2	1:A:328:PHE:CZ	2.52	0.45
6:G:26:ASP:OD1	6:G:27:ARG:N	2.50	0.45
2:B:42:ILE:CG2	2:B:43:ASP:N	2.80	0.45
1:A:15:ARG:NH2	2:B:95:VAL:O	2.42	0.44
3:C:153:ILE:HD11	3:C:177:VAL:HA	1.97	0.44
5:E:60:ALA:HB3	5:E:63:VAL:HG22	1.99	0.44
2:B:158:ASP:OD1	2:B:161:GLN:N	2.47	0.44
5:E:66:ARG:HH22	5:E:86:ARG:HG3	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:G:37:LEU:HD12	6:G:37:LEU:HA	1.87	0.44
1:A:238:PHE:CE2	1:A:254:LEU:HD11	2.51	0.44
5:E:157:ILE:HD11	5:E:214:LEU:HD23	1.99	0.44
2:B:279:THR:HG21	2:B:320:VAL:O	2.18	0.44
2:B:234:ILE:HG22	2:B:250:SER:HB3	1.98	0.44
3:C:155:ILE:HD13	3:C:155:ILE:HA	1.89	0.44
3:C:94:ILE:C	3:C:96:ILE:N	2.71	0.44
3:C:381:ILE:O	3:C:384:ILE:HG12	2.18	0.44
1:A:46:LYS:HB2	1:A:46:LYS:NZ	2.32	0.44
1:A:244:ASN:HB2	1:A:247:LEU:HB2	1.99	0.44
1:A:258:LYS:HG2	1:A:261:LEU:HD13	2.00	0.44
2:B:237:ILE:HB	2:B:248:THR:HB	1.99	0.44
3:C:351:ASN:N	3:C:351:ASN:OD1	2.49	0.44
2:B:165:SER:HB3	2:B:195:LEU:HD12	2.00	0.44
2:B:269:TYR:OH	2:B:304:ALA:O	2.30	0.44
1:A:35:ARG:HG3	1:A:216:VAL:HA	1.99	0.43
3:C:286:LEU:HA	3:C:289:THR:HG22	2.00	0.43
1:A:228:TYR:CZ	1:A:274:GLU:HG3	2.53	0.43
2:B:215:LEU:HD22	2:B:260:LEU:HD22	2.00	0.43
1:A:248:ARG:HG3	1:A:249:THR:HG23	1.99	0.43
2:B:87:TRP:CD1	2:B:87:TRP:N	2.86	0.43
5:E:28:PHE:O	5:E:71:ARG:NH2	2.50	0.43
2:B:144:LEU:HD13	2:B:174:TRP:CD2	2.53	0.43
2:B:251:ASP:HA	2:B:277:GLY:HA3	2.00	0.43
2:B:320:VAL:HA	2:B:336:SER:HB2	2.01	0.43
3:C:365:LEU:HA	3:C:365:LEU:HD23	1.72	0.43
2:B:175:ASP:OD2	2:B:178:THR:OG1	2.33	0.43
2:B:56:LEU:HB3	2:B:87:TRP:CZ3	2.54	0.43
2:B:259:ASP:HB3	2:B:262:ALA:HB3	1.99	0.43
4:D:15:CYS:HA	4:D:18:ASP:HB2	2.01	0.43
5:E:233:LEU:O	5:E:234:GLU:CB	2.67	0.43
2:B:74:LEU:HD23	2:B:74:LEU:HA	1.85	0.43
1:A:8:GLU:N	5:E:173:TYR:HH	2.16	0.42
1:A:273:ILE:HG23	1:A:275:ASP:N	2.34	0.42
5:E:87:SER:HA	5:E:118:VAL:HG21	1.99	0.42
5:E:200:PRO:HG2	5:E:203:PHE:HE2	1.84	0.42
1:A:223:VAL:HB	1:A:256:LEU:HD22	2.01	0.42
2:B:35:LEU:HD23	2:B:266:LEU:HD12	2.01	0.42
5:E:44:LEU:HD23	5:E:44:LEU:HA	1.85	0.42
2:B:272:ASP:N	2:B:272:ASP:OD1	2.51	0.42
3:C:246:ASP:N	3:C:246:ASP:OD1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:343:ARG:HG3	3:C:343:ARG:NH1	2.34	0.42
1:A:221:PHE:HE2	1:A:241:ILE:HD11	1.84	0.42
3:C:139:LEU:HD23	3:C:139:LEU:HA	1.84	0.42
5:E:98:SER:HB2	5:E:106:PRO:HA	2.02	0.42
1:A:272:LYS:O	1:A:273:ILE:HG22	2.18	0.41
3:C:94:ILE:O	3:C:96:ILE:N	2.53	0.41
5:E:36:VAL:HG21	5:E:110:TRP:CZ3	2.55	0.41
2:B:278:ILE:N	2:B:278:ILE:HD12	2.36	0.41
2:B:154:CYS:HB2	2:B:162:ILE:HD11	2.01	0.41
5:E:7:GLY:O	5:E:114:THR:OG1	2.25	0.41
1:A:343:ASP:O	1:A:347:ILE:HG22	2.20	0.41
3:C:157:ILE:H	3:C:157:ILE:HG13	1.74	0.41
1:A:272:LYS:C	1:A:274:GLU:N	2.74	0.41
2:B:256:ARG:HD3	2:B:265:GLU:CD	2.41	0.41
1:A:29:LYS:HA	1:A:32:ARG:HE	1.86	0.41
2:B:274:ILE:HG21	2:B:294:TYR:CE2	2.55	0.41
3:C:189:VAL:HG11	3:C:281:TYR:CE1	2.56	0.41
5:E:157:ILE:HD13	5:E:227:TYR:HD2	1.86	0.41
1:A:36:LEU:HD12	1:A:36:LEU:HA	1.95	0.41
2:B:151:LEU:HD12	2:B:151:LEU:HA	1.90	0.41
3:C:373:ASN:HD22	3:C:373:ASN:HA	1.60	0.41
4:D:18:ASP:OD1	4:D:19:ILE:N	2.54	0.41
5:E:106:PRO:HB2	5:E:175:TYR:HE2	1.86	0.41
5:E:107:PHE:CD2	5:E:230:MET:HE1	2.56	0.40
3:C:160:TYR:HE2	3:C:168:PRO:HD2	1.86	0.40
5:E:169:ASN:OD1	5:E:170:GLY:N	2.54	0.40
2:B:169:THR:HG22	2:B:190:GLY:C	2.41	0.40
4:D:6:LEU:HD23	4:D:6:LEU:HA	1.93	0.40
1:A:206:ARG:HA	1:A:209:TRP:HD1	1.86	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	213/246 (87%)	202 (95%)	10 (5%)	1 (0%)	29	68
2	B	336/377 (89%)	330 (98%)	6 (2%)	0	100	100
3	C	286/603 (47%)	276 (96%)	8 (3%)	2 (1%)	22	61
4	D	19/21 (90%)	18 (95%)	1 (5%)	0	100	100
5	E	228/285 (80%)	224 (98%)	3 (1%)	1 (0%)	34	72
6	G	55/71 (78%)	55 (100%)	0	0	100	100
All	All	1137/1603 (71%)	1105 (97%)	28 (2%)	4 (0%)	38	72

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	96	ILE
1	A	273	ILE
5	E	234	GLU
3	C	95	GLU

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	190/213 (89%)	179 (94%)	11 (6%)	20	53
2	B	279/308 (91%)	270 (97%)	9 (3%)	39	69
3	C	227/536 (42%)	209 (92%)	18 (8%)	12	41
4	D	19/21 (90%)	15 (79%)	4 (21%)	1	5
5	E	193/228 (85%)	185 (96%)	8 (4%)	30	63
6	G	45/58 (78%)	44 (98%)	1 (2%)	52	78
All	All	953/1364 (70%)	902 (95%)	51 (5%)	26	55

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

Mol	Chain	Res	Type
1	A	24	ARG

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Mol	Chain	Res	Type
1	A	35	ARG
1	A	183	PHE
1	A	194	PHE
1	A	227	ASP
1	A	239	LYS
1	A	273	ILE
1	A	277	PHE
1	A	309	GLU
1	A	321	ARG
1	A	333	ASP
2	B	9	LEU
2	B	50	MET
2	B	115	ASN
2	B	239	PHE
2	B	261	ARG
2	B	263	ASP
2	B	319	ARG
2	B	322	CYS
2	B	336	SER
3	C	94	ILE
3	C	96	ILE
3	C	100	PHE
3	C	129	ASN
3	C	131	CYS
3	C	150	HIS
3	C	181	GLN
3	C	184	SER
3	C	205	SER
3	C	206	TRP
3	C	207	SER
3	C	240	PHE
3	C	278	PHE
3	C	280	PHE
3	C	282	PHE
3	C	284	LEU
3	C	293	TYR
3	C	373	ASN
4	D	13	TYR
4	D	14	PHE
4	D	15	CYS
4	D	17	LEU
5	E	21	CYS

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Mol	Chain	Res	Type
5	E	28	PHE
5	E	42	LYS
5	E	75	LYS
5	E	97	ARG
5	E	190	TYR
5	E	206	SER
5	E	247	LEU
6	G	44	HIS

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

Mol	Chain	Res	Type
3	C	181	GLN
5	E	231	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 monosaccharides 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.

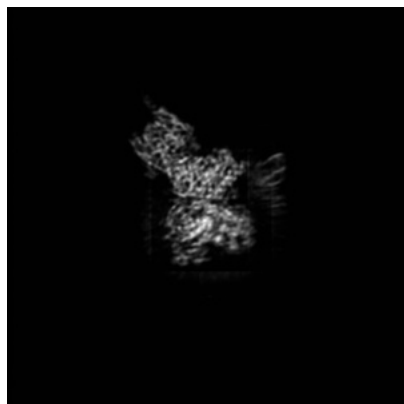
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34667. 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

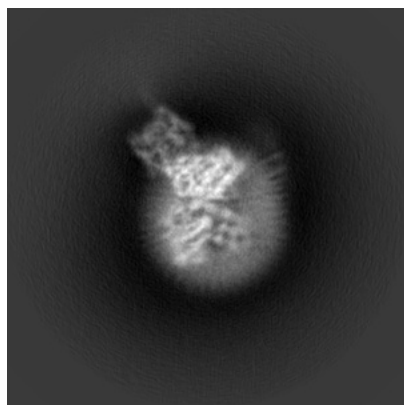


Y

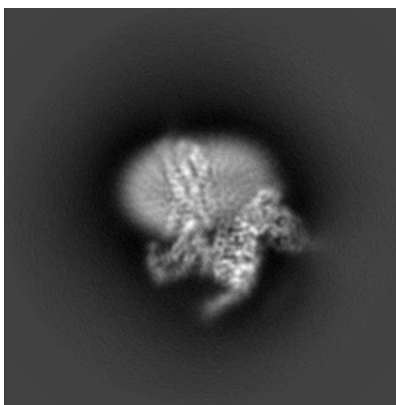


Z

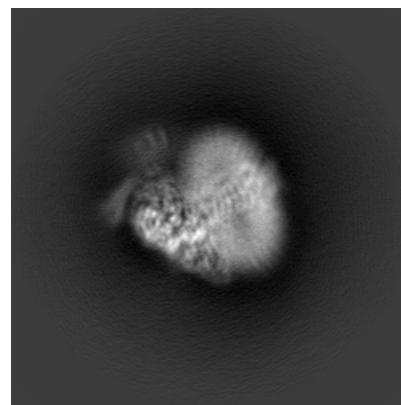
6.1.2 Raw map



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: 128

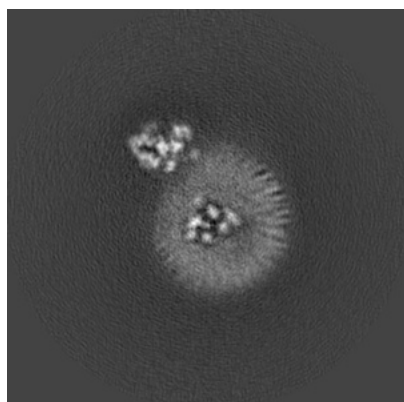


Y Index: 128

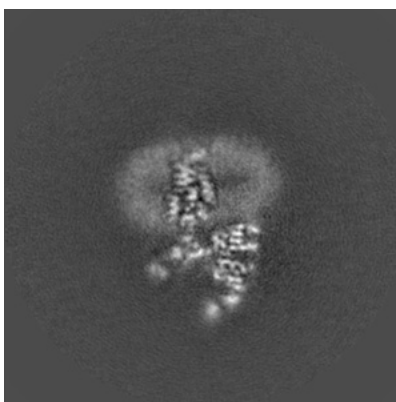


Z Index: 128

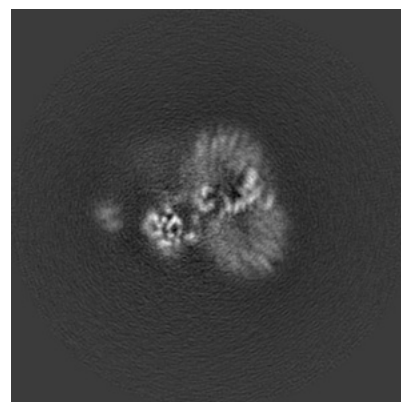
6.2.2 Raw map



X Index: 128



Y Index: 128



Z Index: 128

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: 102

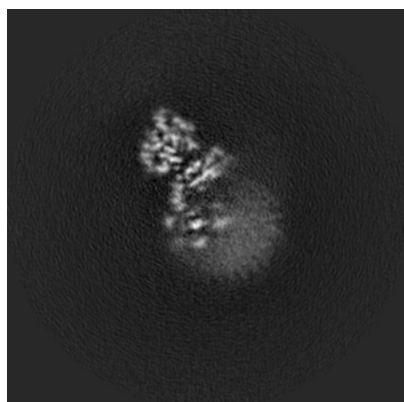


Y Index: 117

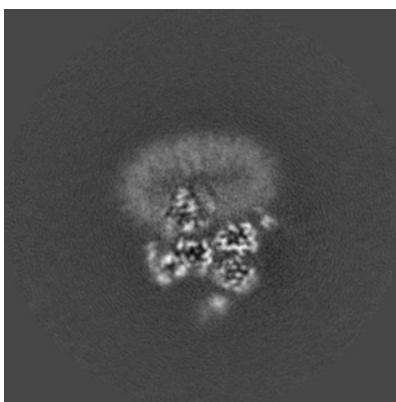


Z Index: 149

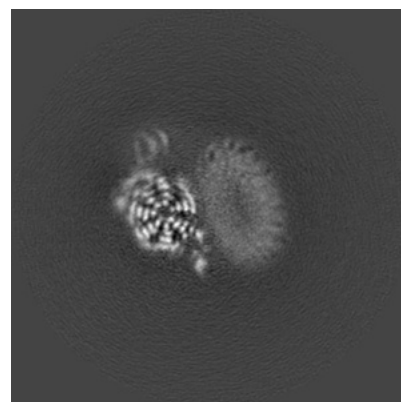
6.3.2 Raw map



X Index: 116



Y Index: 121



Z Index: 150

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

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

6.4.1 Primary map



X

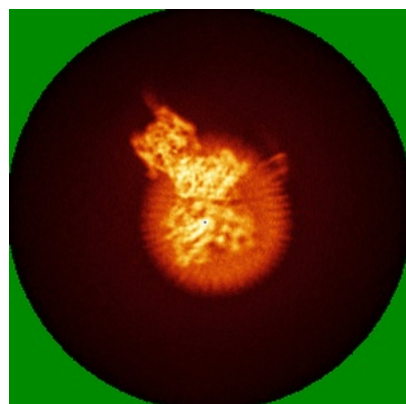


Y



Z

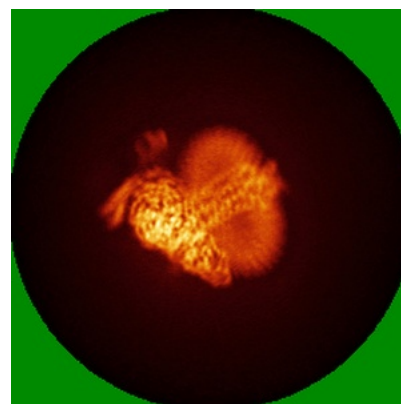
6.4.2 Raw map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.013. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

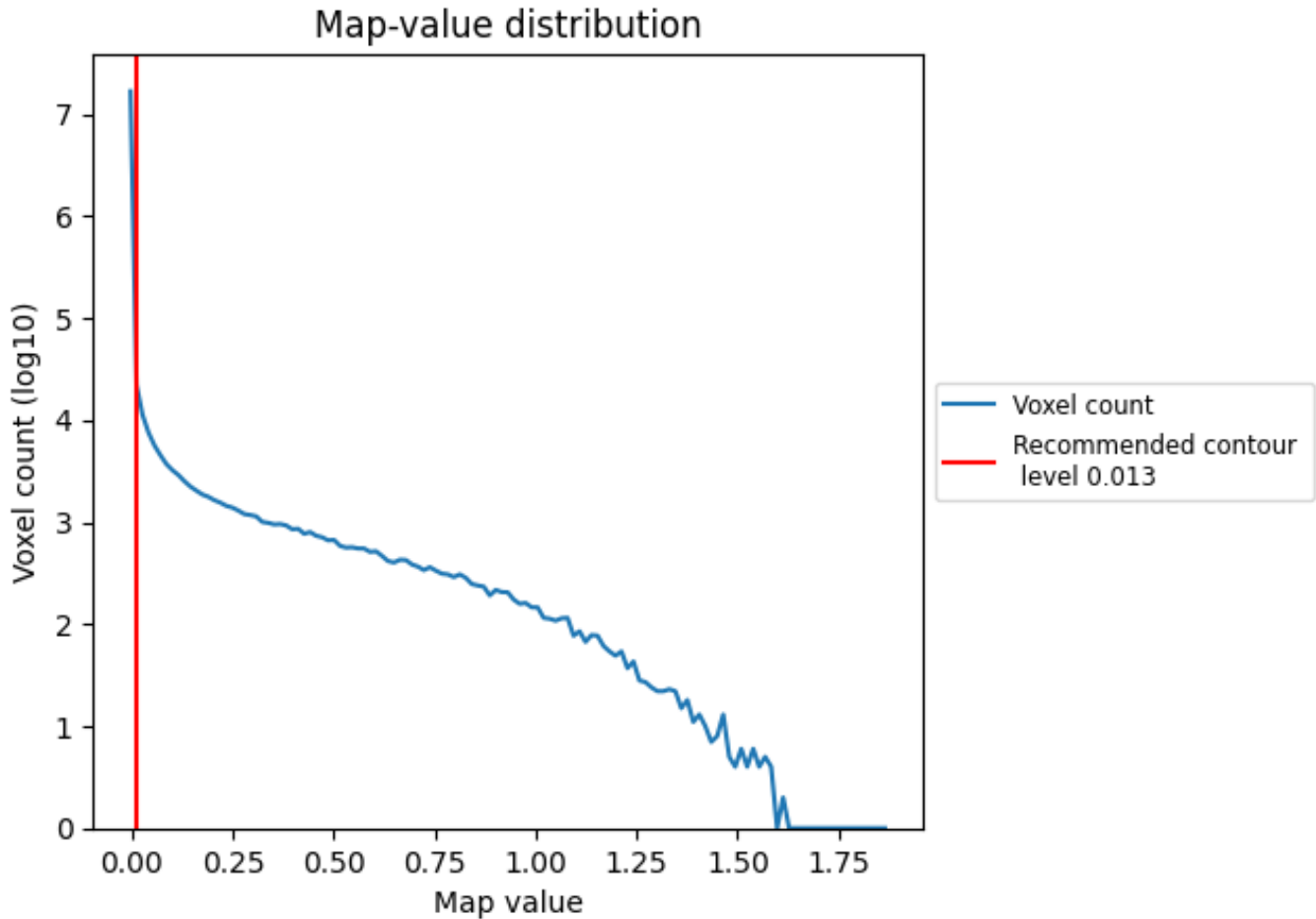
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

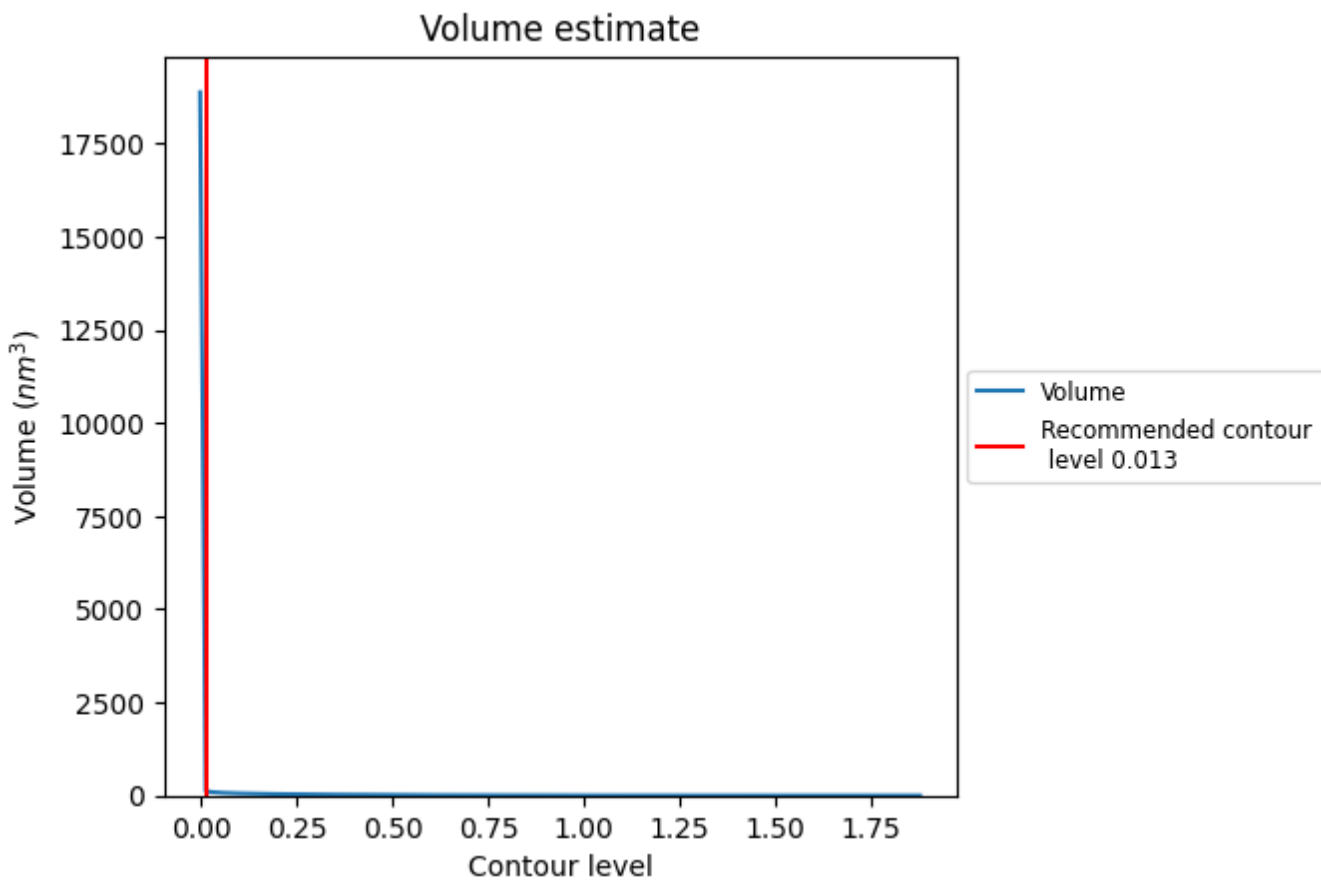
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

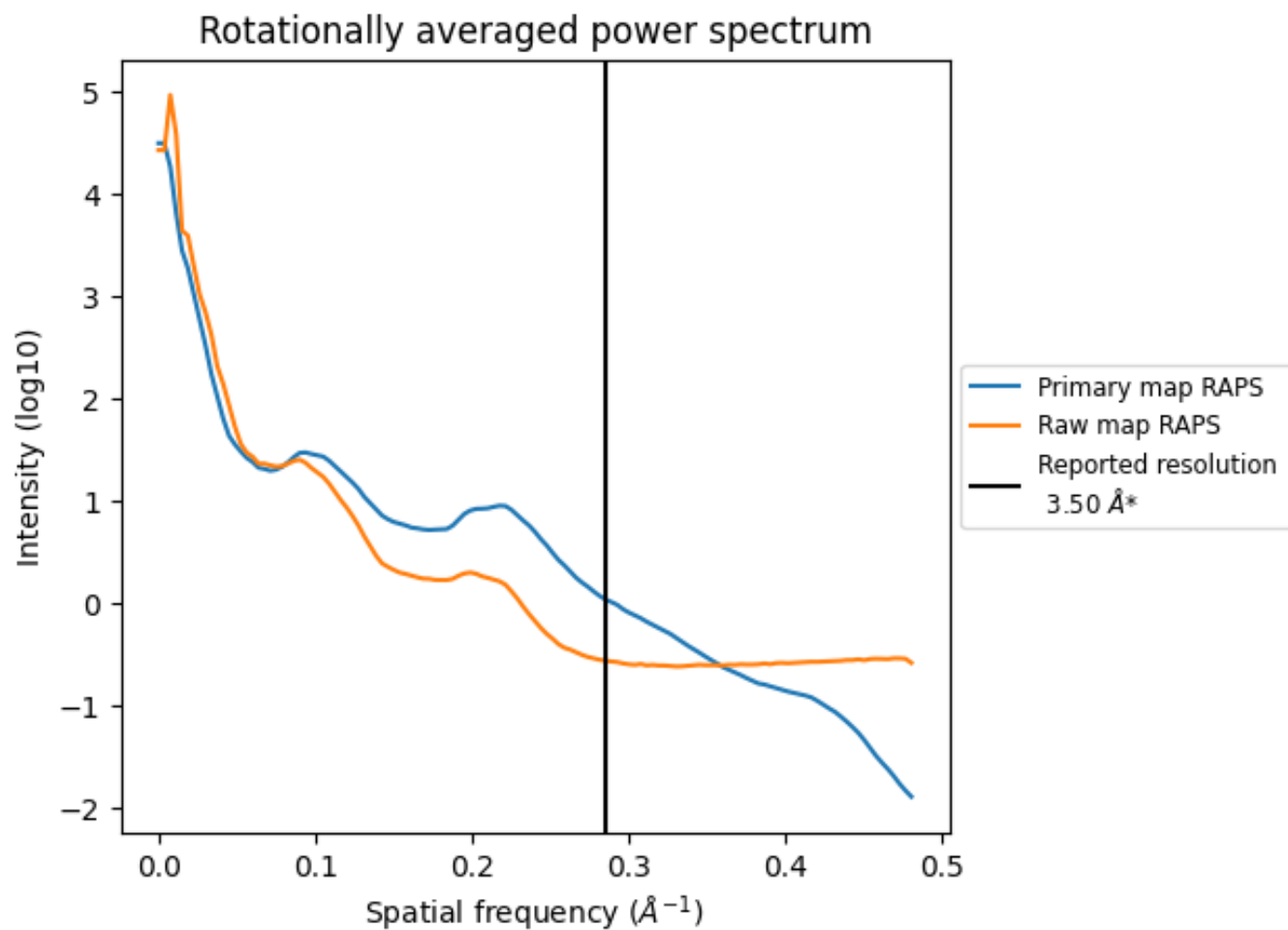
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 117 nm^3 ; this corresponds to an approximate mass of 105 kDa.

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

7.3 Rotationally averaged power spectrum i

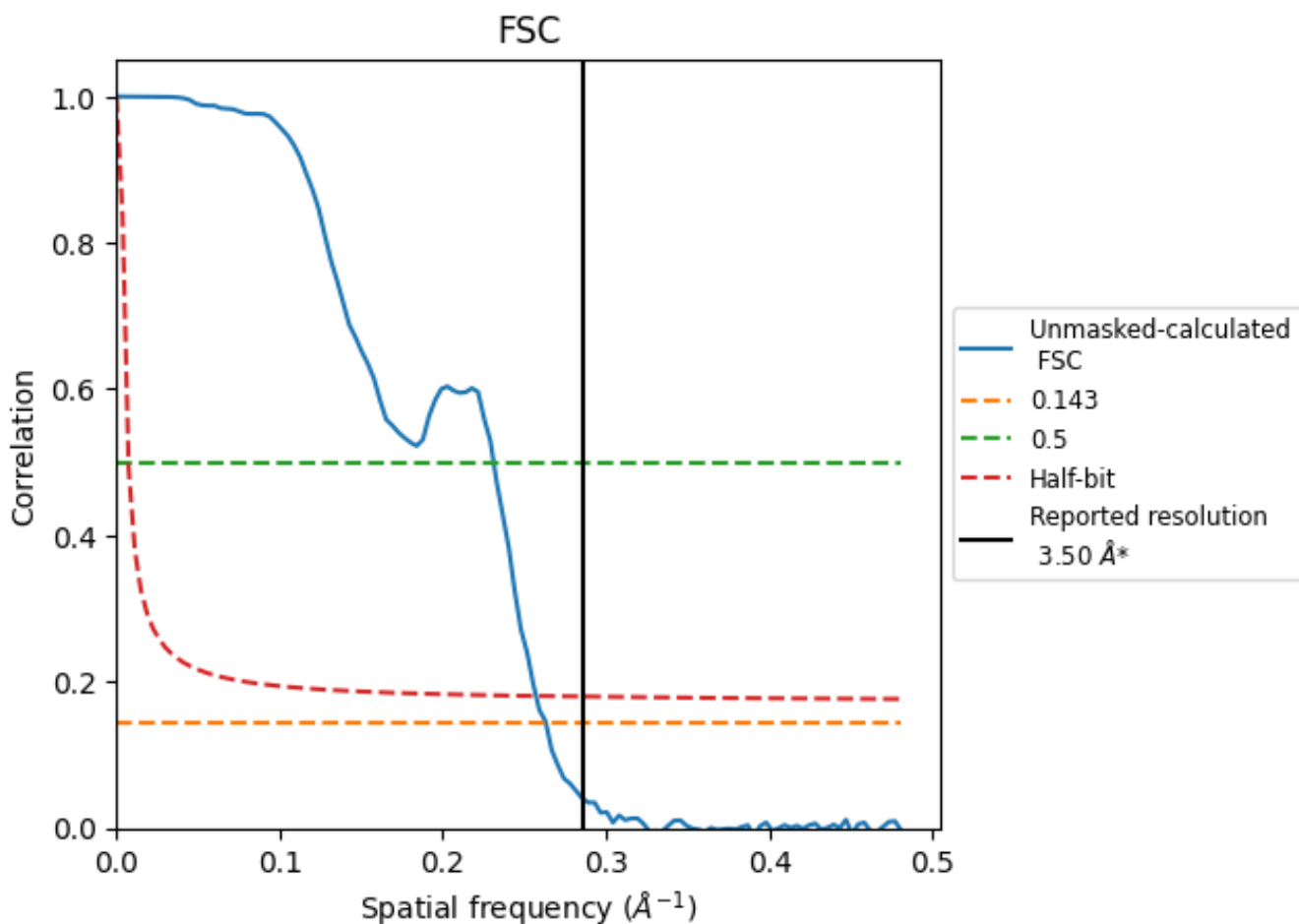


*Reported resolution corresponds to spatial frequency of 0.286 \AA^{-1}

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



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

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.80	4.33	3.89

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

9 Map-model fit [i](#)

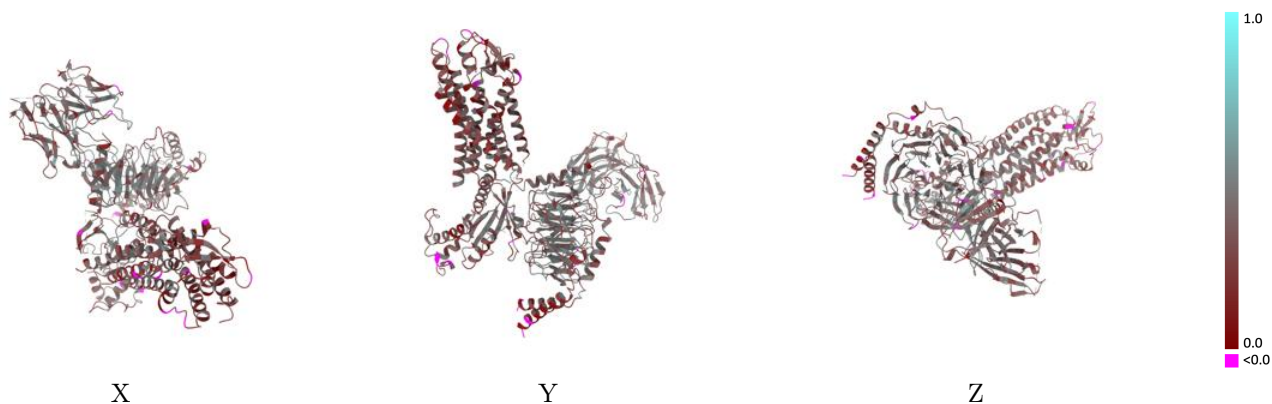
This section contains information regarding the fit between EMDB map EMD-34667 and PDB model 8HCX. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)



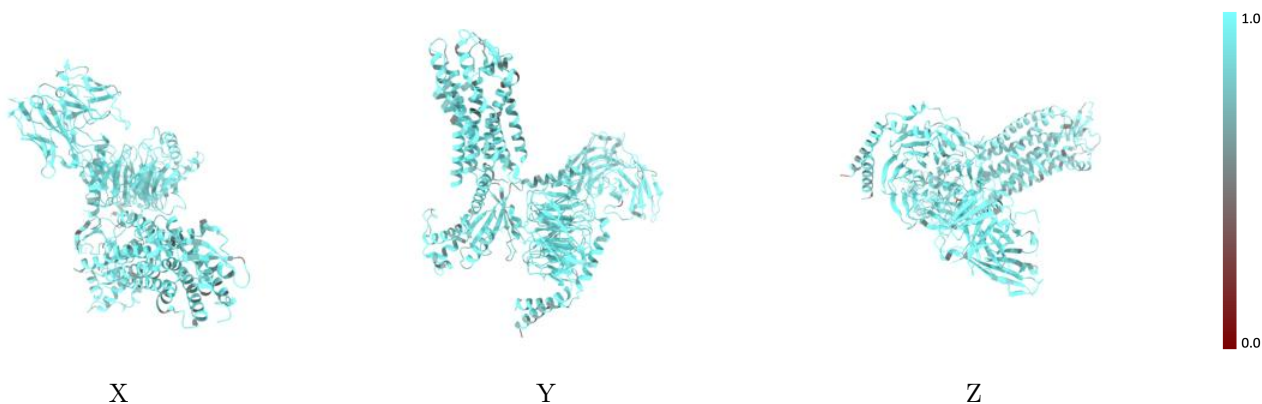
The images above show the 3D surface view of the map at the recommended contour level 0.013 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



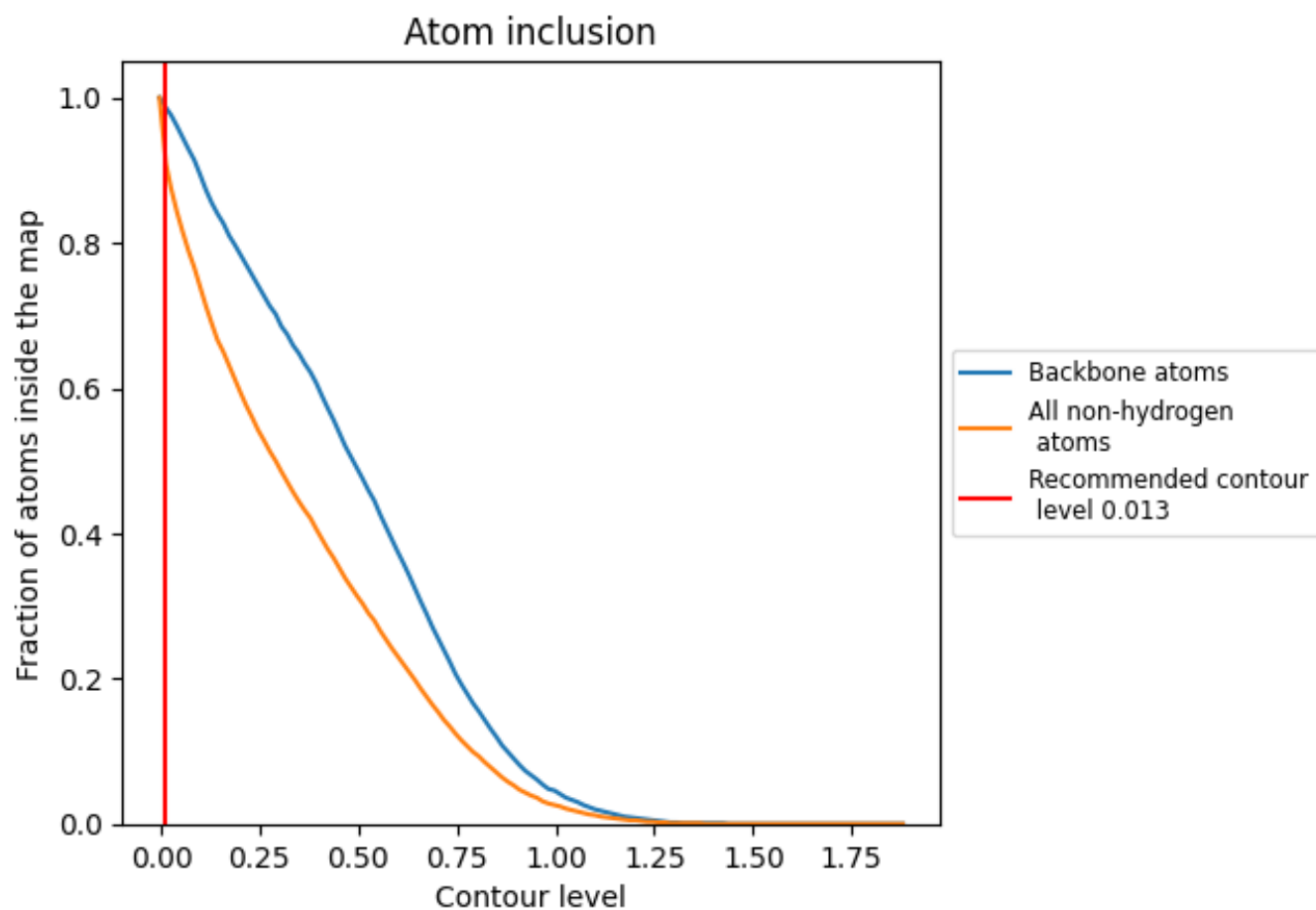
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.013).















9.4 Atom inclusion [i](#)



At the recommended contour level, 99% of all backbone atoms, 91% 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 (0.013) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9090	 0.3460
A	 0.8940	 0.3300
B	 0.9390	 0.3910
C	 0.8810	 0.2920
D	 0.8520	 0.2520
E	 0.9290	 0.3820
G	 0.8800	 0.3090

