Full wwPDB EM Validation Report

Apr 24, 2025 – 04:21 PM JST

PDB ID	:	9IK8 / pdb_00009ik8
EMDB ID	:	EMD-60650
Title	:	Cryo-EM Structure of SSTR1-Gi SST analogs complex
Authors	:	Wong, T.S.; Zeng, Z.C.; Xiong, T.T.; Gan, S.Y.; Du, Y.
Deposited on	:	2024-06-26
Resolution	:	2.82 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:

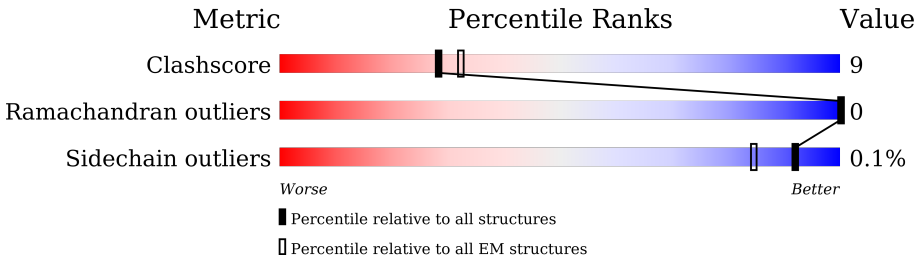
EMDB validation analysis	:	0.0.1.dev117
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.42

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 2.82 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	D	393	12% 59% 11% 30%
2	E	338	57% 11% 31%
3	B	373	79% 11% 10%
4	C	71	6% 73% 25%
5	A	350	53% 9% 39%
6	F	6	33% 67%

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
6	TY5	F	3	-	-	X	-
6	004	F	6	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16815 atoms, of which 8318 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 Somatostatin receptor type 1.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	D	274	Total	C	H	N	O	S	0	0
			4358	1409	2237	348	342	22		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	392	LEU	-	expression tag	UNP P30872
D	393	GLU	-	expression tag	UNP P30872

- Molecule 2 is a protein called scFv16.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	E	232	Total	C	H	N	O	S	0	0
			3416	1116	1668	295	328	9		

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	B	336	Total	C	H	N	O	S	0	0
			4912	1563	2389	450	491	19		

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-21	MET	-	initiating methionine	UNP P62873
B	-20	HIS	-	expression tag	UNP P62873
B	-19	HIS	-	expression tag	UNP P62873
B	-18	HIS	-	expression tag	UNP P62873
B	-17	HIS	-	expression tag	UNP P62873
B	-16	HIS	-	expression tag	UNP P62873
B	-15	HIS	-	expression tag	UNP P62873
B	-14	HIS	-	expression tag	UNP P62873

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	HIS	-	expression tag	UNP P62873
B	-12	HIS	-	expression tag	UNP P62873
B	-11	HIS	-	expression tag	UNP P62873
B	-10	LEU	-	expression tag	UNP P62873
B	-9	GLU	-	expression tag	UNP P62873
B	-8	VAL	-	expression tag	UNP P62873
B	-7	LEU	-	expression tag	UNP P62873
B	-6	PHE	-	expression tag	UNP P62873
B	-5	GLN	-	expression tag	UNP P62873
B	-4	GLY	-	expression tag	UNP P62873
B	-3	PRO	-	expression tag	UNP P62873
B	-2	GLY	-	expression tag	UNP P62873
B	-1	SER	-	expression tag	UNP P62873
B	0	SER	-	expression tag	UNP P62873
B	1	GLY	-	expression tag	UNP P62873
B	341	VAL	-	expression tag	UNP P62873
B	342	SER	-	expression tag	UNP P62873
B	343	GLY	-	expression tag	UNP P62873
B	344	TRP	-	expression tag	UNP P62873
B	345	ARG	-	expression tag	UNP P62873
B	346	LEU	-	expression tag	UNP P62873
B	347	PHE	-	expression tag	UNP P62873
B	348	LYS	-	expression tag	UNP P62873
B	349	LYS	-	expression tag	UNP P62873
B	350	ILE	-	expression tag	UNP P62873
B	351	SER	-	expression tag	UNP P62873

- Molecule 4 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	C	53	Total	C	H	N	O	S	0	0
			749	241	372	66	67	3		

- Molecule 5 is a protein called Guanine nucleotide-binding protein G(i) subunit alpha-1.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	A	215	Total	C	H	N	O	S	0	0
			3237	1058	1586	284	297	12		

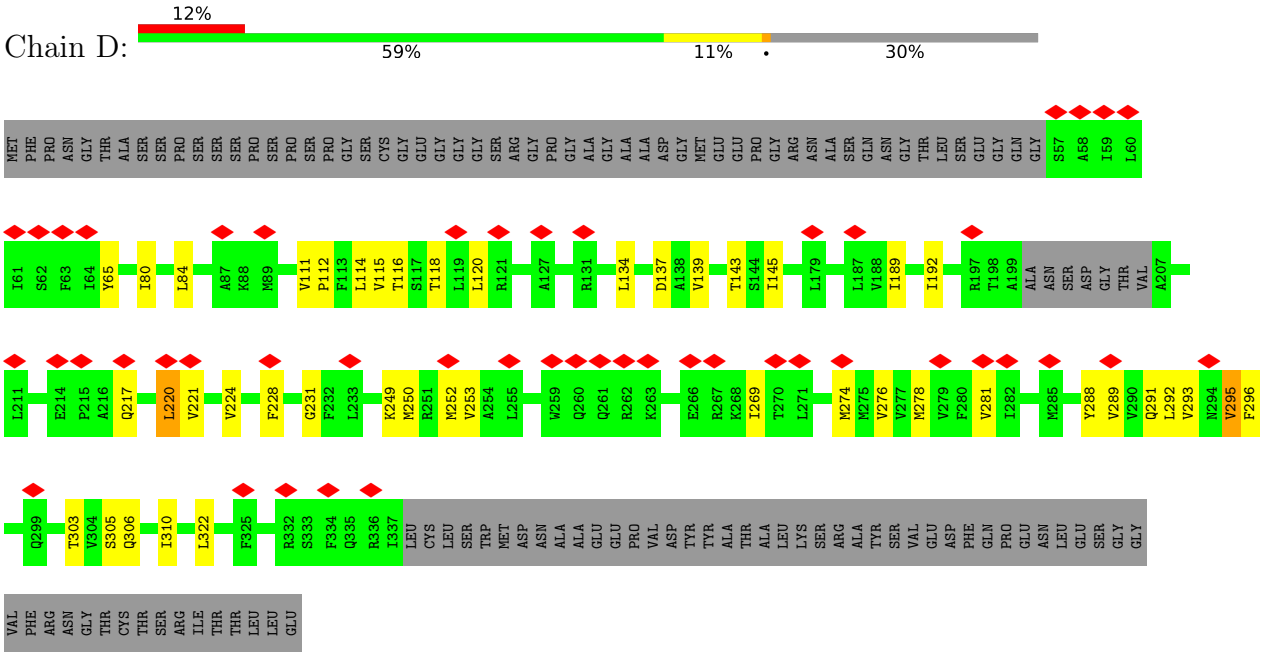
- Molecule 6 is a protein called DTR-LYS-TY5-PHA-A1D5E-004.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	6	Total	C	H	N	O	0	0
			143	58	66	10	9		

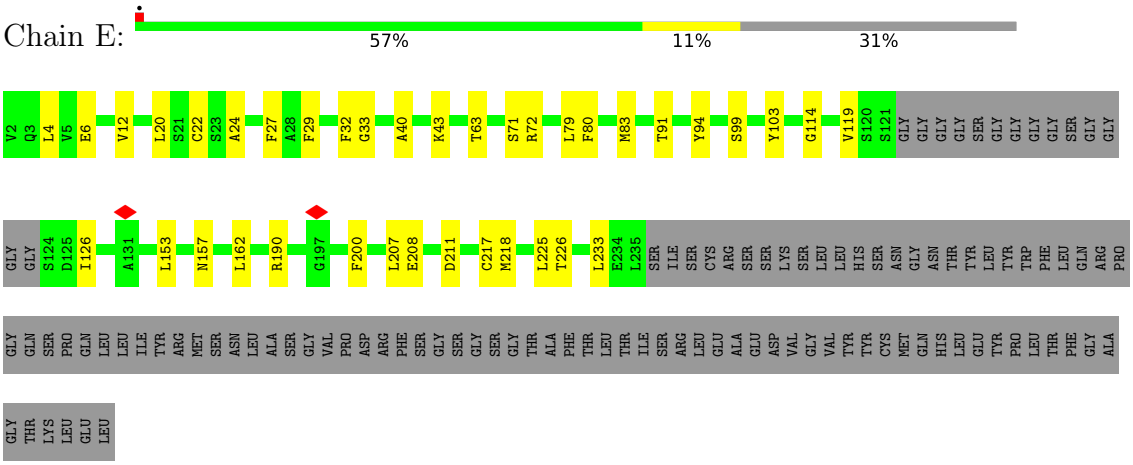
3 Residue-property plots (i)

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: Somatostatin receptor type 1

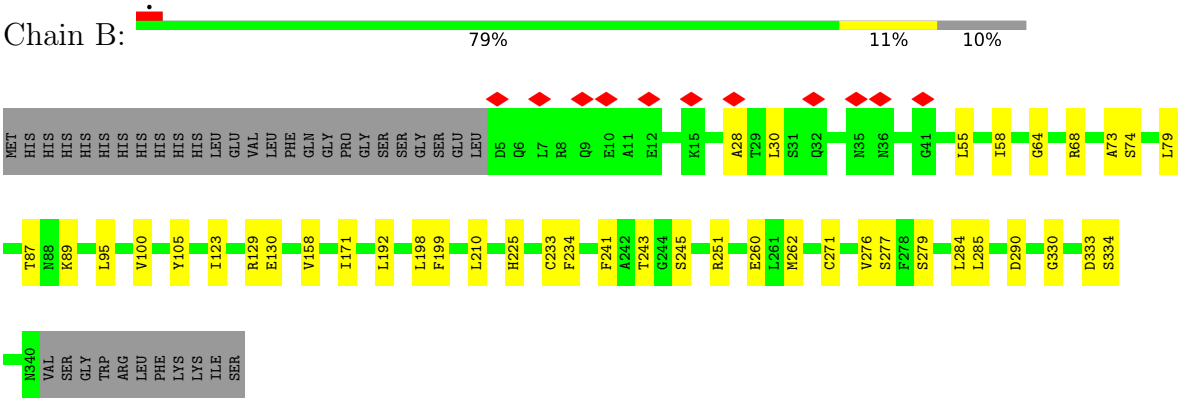


• Molecule 2: scFv16



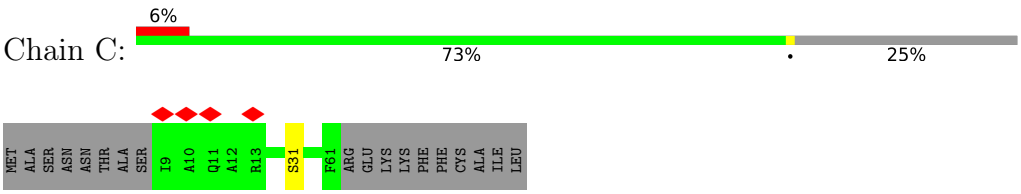
• Molecule 3: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1

Chain B:



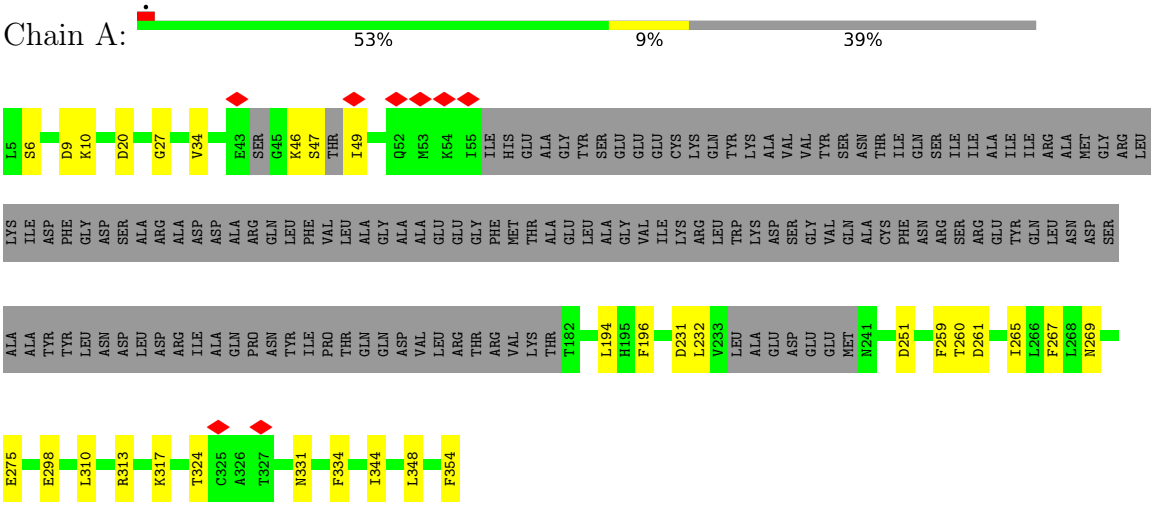
• Molecule 4: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2

Chain C:



• Molecule 5: Guanine nucleotide-binding protein G(i) subunit alpha-1

Chain A:



• Molecule 6: DTR-LYS-TY5-PHA-A1D5E-004

Chain F:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	722906	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 ($e^-/Å^2$)	53.5	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	1400	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOCONTINUUM (6k x 4k)	Depositor
Maximum map value	3.181	Depositor
Minimum map value	-2.222	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.037	Depositor
Recommended contour level	0.12	Depositor
Map size (Å)	318.72, 318.72, 318.72	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PHA, DTR, 004, A1D5E, TY5

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	D	0.31	0/2165	0.54	1/2951 (0.0%)
2	E	0.32	0/1792	0.55	0/2434
3	B	0.39	0/2570	0.61	0/3493
4	C	0.38	0/383	0.49	0/522
5	A	0.42	0/1678	0.54	0/2255
6	F	1.28	0/8	1.34	0/8
All	All	0.36	0/8596	0.56	1/11663 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	295	VAL	CB-CA-C	-5.71	100.55	111.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	D	2121	2237	2246	72	0
2	E	1748	1668	1669	26	0
3	B	2523	2389	2389	35	0
4	C	377	372	372	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	1651	1586	1584	31	0
6	F	77	66	53	38	0
All	All	8497	8318	8313	156	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 9.

All (156) 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:D:114:LEU:HD12	6:F:3:TY5:C54	1.25	1.58
1:D:114:LEU:CD1	6:F:3:TY5:H54	1.42	1.46
1:D:295:VAL:CG1	6:F:6:004:HD2	1.47	1.42
1:D:114:LEU:HD12	6:F:3:TY5:C55	1.56	1.34
1:D:269:ILE:HD11	5:A:354:PHE:CD1	1.66	1.31
1:D:295:VAL:CG1	6:F:6:004:CD2	2.13	1.27
1:D:114:LEU:CD1	6:F:3:TY5:C54	2.06	1.21
1:D:114:LEU:HD13	6:F:3:TY5:H54	1.30	1.10
1:D:220:LEU:HD13	6:F:6:004:CE	1.89	1.01
1:D:269:ILE:CD1	5:A:354:PHE:HD1	1.74	1.01
1:D:269:ILE:HD11	5:A:354:PHE:HD1	0.81	0.97
1:D:295:VAL:HG12	6:F:6:004:CD2	1.94	0.96
1:D:295:VAL:HG13	6:F:6:004:CD2	1.84	0.93
1:D:295:VAL:HG13	6:F:6:004:HD2	0.93	0.92
1:D:289:VAL:O	1:D:293:VAL:HG23	1.71	0.90
1:D:137:ASP:OD2	6:F:3:TY5:C52	2.21	0.88
1:D:224:VAL:HG21	6:F:6:004:CE	2.04	0.86
1:D:114:LEU:HD12	6:F:3:TY5:H55	1.59	0.84
5:A:49:ILE:HD11	5:A:269:ASN:HD22	1.42	0.83
3:B:198:LEU:HD13	3:B:210:LEU:HD11	1.65	0.77
1:D:253:VAL:HG21	5:A:344:ILE:HD13	1.66	0.76
1:D:114:LEU:CD1	6:F:3:TY5:C55	2.52	0.75
1:D:291:GLN:NE2	6:F:1:DTR:O	2.17	0.74
2:E:29:PHE:O	2:E:72:ARG:NH2	2.20	0.73
1:D:220:LEU:HD13	6:F:6:004:CD1	2.19	0.72
5:A:46:LYS:O	5:A:47:SER:OG	2.07	0.71
1:D:220:LEU:HD12	1:D:221:VAL:N	2.04	0.71
3:B:192:LEU:HD23	3:B:199:PHE:HB3	1.73	0.70
5:A:260:THR:HG22	5:A:313:ARG:NH1	2.06	0.70
3:B:30:LEU:HD22	3:B:262:MET:HE1	1.74	0.69
1:D:114:LEU:HB2	6:F:3:TY5:C54	2.24	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:333:ASP:O	3:B:334:SER:OG	2.11	0.67
2:E:32:PHE:O	2:E:72:ARG:NH2	2.26	0.66
2:E:20:LEU:HD13	2:E:83:MET:HE3	1.77	0.66
5:A:251:ASP:HB2	5:A:310:LEU:HD23	1.79	0.65
1:D:295:VAL:CG1	6:F:6:004:CE	2.75	0.64
1:D:80:ILE:HG22	1:D:84:LEU:HD11	1.78	0.64
1:D:253:VAL:HG21	5:A:344:ILE:CD1	2.26	0.64
2:E:6:GLU:OE1	2:E:114:GLY:N	2.30	0.64
3:B:123:ILE:HD13	3:B:171:ILE:HD12	1.80	0.64
3:B:95:LEU:HD13	3:B:100:VAL:HG21	1.80	0.63
3:B:225:HIS:CE1	3:B:243:THR:HG1	2.16	0.63
3:B:30:LEU:HD22	3:B:262:MET:CE	2.30	0.62
1:D:295:VAL:HG23	1:D:296:PHE:N	2.14	0.62
1:D:114:LEU:HG	6:F:3:TY5:H44	1.82	0.62
5:A:260:THR:HG22	5:A:313:ARG:HH11	1.64	0.62
1:D:112:PRO:O	1:D:116:THR:HG23	2.00	0.61
2:E:20:LEU:HD13	2:E:83:MET:CE	2.31	0.60
3:B:58:ILE:HD11	3:B:330:GLY:CA	2.31	0.60
1:D:114:LEU:HG	6:F:3:TY5:CE2	2.32	0.60
3:B:276:VAL:HG13	3:B:285:LEU:HD11	1.83	0.60
1:D:228:PHE:HD1	1:D:288:TYR:HB3	1.65	0.59
2:E:103:TYR:CZ	3:B:68:ARG:HD3	2.37	0.59
3:B:225:HIS:ND1	3:B:245:SER:OG	2.33	0.59
3:B:87:THR:O	3:B:87:THR:HG22	2.02	0.59
3:B:225:HIS:NE2	3:B:243:THR:OG1	2.31	0.59
3:B:95:LEU:CD1	3:B:100:VAL:HG21	2.33	0.58
1:D:217:GLN:HA	1:D:220:LEU:HG	1.86	0.58
1:D:249:LYS:O	1:D:253:VAL:HG22	2.05	0.57
1:D:295:VAL:CG2	1:D:296:PHE:N	2.68	0.56
5:A:231:ASP:OD2	5:A:232:LEU:N	2.38	0.56
1:D:252:MET:HG3	1:D:253:VAL:HG13	1.86	0.56
2:E:233:LEU:HD23	2:E:233:LEU:O	2.04	0.56
1:D:220:LEU:HB2	6:F:6:004:HD1	1.86	0.56
5:A:259:PHE:O	5:A:317:LYS:NZ	2.36	0.56
2:E:33:GLY:N	2:E:99:SER:O	2.39	0.55
2:E:40:ALA:HB3	2:E:43:LYS:CG	2.37	0.55
2:E:218:MET:HE1	2:E:225:LEU:HD13	1.88	0.55
2:E:153:LEU:HD12	2:E:200:PHE:CE2	2.42	0.54
1:D:220:LEU:HD12	1:D:220:LEU:C	2.28	0.54
1:D:305:SER:OG	6:F:4:PHA:HB3	2.08	0.54
3:B:271:CYS:HB3	3:B:290:ASP:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:114:LEU:HB2	6:F:3:TY5:C53	2.38	0.54
1:D:250:MET:SD	5:A:348:LEU:HD11	2.48	0.54
1:D:274:MET:O	1:D:274:MET:HG2	2.08	0.54
1:D:224:VAL:CG2	6:F:6:004:CE	2.82	0.53
1:D:228:PHE:CD1	1:D:288:TYR:HB3	2.43	0.53
5:A:49:ILE:HD13	5:A:324:THR:HG21	1.89	0.53
5:A:348:LEU:HB3	5:A:354:PHE:HB3	1.90	0.53
5:A:354:PHE:CD2	5:A:354:PHE:OXT	2.61	0.53
1:D:303:THR:HG23	1:D:306:GLN:NE2	2.23	0.53
3:B:89:LYS:NZ	5:A:20:ASP:OD1	2.34	0.53
5:A:34:VAL:O	5:A:34:VAL:HG23	2.09	0.53
3:B:158:VAL:HG11	3:B:192:LEU:HD21	1.90	0.52
5:A:261:ASP:OD1	5:A:261:ASP:N	2.41	0.52
2:E:4:LEU:HD22	2:E:22:CYS:SG	2.49	0.52
1:D:65:TYR:CE1	1:D:310:ILE:HG22	2.45	0.52
3:B:64:GLY:HA2	3:B:105:TYR:CD2	2.45	0.52
1:D:137:ASP:OD2	6:F:3:TY5:C53	2.58	0.52
1:D:139:VAL:O	1:D:143:THR:HG22	2.10	0.52
3:B:129:ARG:O	3:B:130:GLU:CB	2.57	0.52
2:E:103:TYR:CE1	3:B:68:ARG:NH1	2.78	0.51
1:D:220:LEU:CD1	6:F:6:004:CE	2.76	0.51
1:D:269:ILE:CD1	5:A:354:PHE:CD1	2.62	0.51
3:B:58:ILE:HD11	3:B:330:GLY:HA3	1.92	0.51
1:D:145:ILE:HD13	1:D:231:GLY:O	2.10	0.51
2:E:40:ALA:HB3	2:E:43:LYS:HG2	1.92	0.51
3:B:198:LEU:HD12	3:B:198:LEU:O	2.09	0.51
3:B:279:SER:OG	3:B:284:LEU:HB2	2.11	0.51
5:A:298:GLU:OE2	5:A:298:GLU:HA	2.10	0.51
2:E:94:TYR:O	2:E:114:GLY:HA2	2.11	0.51
3:B:73:ALA:HB2	3:B:79:LEU:HD12	1.93	0.51
2:E:24:ALA:HB1	2:E:27:PHE:CZ	2.47	0.50
2:E:190:ARG:NH2	2:E:211:ASP:OD2	2.43	0.50
1:D:217:GLN:O	1:D:220:LEU:HG	2.10	0.50
1:D:137:ASP:OD2	6:F:3:TY5:H52	2.07	0.50
1:D:134:LEU:HD13	6:F:3:TY5:H49	1.93	0.50
2:E:71:SER:HB2	2:E:80:PHE:HB2	1.93	0.50
2:E:162:LEU:HD11	2:E:217:CYS:SG	2.52	0.50
2:E:12:VAL:HG23	2:E:119:VAL:HG12	1.93	0.50
2:E:207:LEU:HD23	2:E:208:GLU:N	2.27	0.50
3:B:55:LEU:HD13	5:A:27:GLY:HA3	1.92	0.49
3:B:251:ARG:NH1	3:B:260:GLU:OE1	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:265:ILE:HG23	5:A:334:PHE:HE1	1.77	0.49
2:E:63:THR:O	2:E:63:THR:HG22	2.12	0.49
3:B:225:HIS:CE1	3:B:251:ARG:HD2	2.47	0.49
1:D:80:ILE:HG22	1:D:84:LEU:CD1	2.41	0.49
3:B:123:ILE:HD13	3:B:171:ILE:CD1	2.43	0.49
1:D:276:VAL:HG22	1:D:322:LEU:HD13	1.95	0.49
1:D:295:VAL:HG11	6:F:6:004:CD2	2.30	0.49
2:E:157:ASN:ND2	5:A:9:ASP:OD2	2.46	0.48
2:E:91:THR:OG1	2:E:119:VAL:HG22	2.13	0.48
5:A:275:GLU:HA	5:A:275:GLU:OE2	2.13	0.48
1:D:221:VAL:HG12	1:D:295:VAL:HG21	1.95	0.48
2:E:126:ILE:O	2:E:226:THR:HG21	2.14	0.48
5:A:6:SER:O	5:A:10:LYS:HG3	2.13	0.48
1:D:114:LEU:O	1:D:118:THR:HG23	2.14	0.48
3:B:233:CYS:O	3:B:241:PHE:HB2	2.14	0.47
1:D:295:VAL:HG11	6:F:6:004:HD2	1.75	0.47
5:A:324:THR:HG23	5:A:331:ASN:OD1	2.14	0.47
6:F:4:PHA:O	6:F:6:004:HA	2.14	0.47
1:D:111:VAL:O	1:D:115:VAL:HG23	2.15	0.47
3:B:58:ILE:HG22	3:B:74:SER:HB2	1.96	0.47
1:D:217:GLN:O	1:D:221:VAL:HG22	2.15	0.47
1:D:278:MET:HA	1:D:281:VAL:HG12	1.97	0.47
6:F:3:TY5:CD1	6:F:3:TY5:N	2.78	0.46
2:E:22:CYS:HB3	2:E:79:LEU:HB3	1.98	0.46
3:B:73:ALA:HB2	3:B:79:LEU:CD1	2.45	0.46
5:A:49:ILE:HD12	5:A:267:PHE:HD2	1.81	0.45
5:A:194:LEU:HD23	5:A:196:PHE:CZ	2.51	0.45
3:B:277:SER:O	3:B:285:LEU:HD12	2.16	0.45
1:D:295:VAL:HG11	6:F:6:004:CE	2.46	0.45
1:D:111:VAL:N	1:D:112:PRO:HD2	2.33	0.44
1:D:292:LEU:O	1:D:293:VAL:C	2.54	0.43
1:D:114:LEU:HG	6:F:3:TY5:CD2	2.49	0.43
1:D:134:LEU:HD23	1:D:192:ILE:HG23	2.01	0.42
6:F:4:PHA:CD2	6:F:4:PHA:C	2.97	0.42
1:D:116:THR:O	1:D:120:LEU:HB2	2.18	0.42
5:A:354:PHE:OXT	5:A:354:PHE:CG	2.73	0.42
3:B:192:LEU:O	3:B:234:PHE:CE2	2.73	0.42
6:F:1:DTR:CD1	6:F:2:LYS:HG2	2.49	0.42
1:D:289:VAL:O	1:D:293:VAL:CG2	2.56	0.41
1:D:114:LEU:CD1	6:F:3:TY5:H55	2.37	0.41
1:D:189:ILE:O	1:D:192:ILE:HD12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:348:LEU:HB3	5:A:354:PHE:CB	2.51	0.41
3:B:28:ALA:HB2	4:C:31:SER:OG	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	D	270/393 (69%)	250 (93%)	20 (7%)	0	100	100
2	E	228/338 (68%)	220 (96%)	8 (4%)	0	100	100
3	B	334/373 (90%)	317 (95%)	17 (5%)	0	100	100
4	C	51/71 (72%)	51 (100%)	0	0	100	100
5	A	205/350 (59%)	201 (98%)	4 (2%)	0	100	100
6	F	1/6 (17%)	1 (100%)	0	0	100	100
All	All	1089/1531 (71%)	1040 (96%)	49 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	D	232/332 (70%)	231 (100%)	1 (0%)	89	96

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	184/278 (66%)	184 (100%)	0	100	100
3	B	263/312 (84%)	263 (100%)	0	100	100
4	C	35/58 (60%)	35 (100%)	0	100	100
5	A	165/302 (55%)	165 (100%)	0	100	100
6	F	1/1 (100%)	1 (100%)	0	100	100
All	All	880/1283 (69%)	879 (100%)	1 (0%)	92	98

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	220	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	306	GLN
4	C	59	ASN
5	A	269	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

5 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
6	TY5	F	3	6	19,20,21	1.35	3 (15%)	22,25,27	0.91	1 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
6	A1D5E	F	5	6	11,14,15	4.11	5 (45%)	11,17,19	2.44	4 (36%)
6	004	F	6	6	9,10,11	1.07	0	9,12,14	1.79	2 (22%)
6	DTR	F	1	6	13,15,16	1.51	2 (15%)	13,20,22	0.97	1 (7%)
6	PHA	F	4	6	10,11,11	1.06	1 (10%)	10,13,13	0.62	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	TY5	F	3	6	-	6/10/11/13	0/2/2/2
6	A1D5E	F	5	6	-	5/8/19/21	0/1/1/1
6	004	F	6	6	-	1/4/6/8	0/1/1/1
6	DTR	F	1	6	-	1/4/6/8	0/2/2/2
6	PHA	F	4	6	-	0/5/6/6	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	5	A1D5E	CB-CG	-8.57	1.34	1.52
6	F	5	A1D5E	C72-N74	7.05	1.49	1.34
6	F	5	A1D5E	CB-CA	6.21	1.68	1.54
6	F	5	A1D5E	CD-N	3.46	1.59	1.47
6	F	5	A1D5E	O71-C72	3.11	1.40	1.35
6	F	3	TY5	CE2-CD2	2.99	1.44	1.38
6	F	3	TY5	CD1-CE1	2.76	1.43	1.38
6	F	1	DTR	CZ3-CE3	2.60	1.42	1.36
6	F	3	TY5	OH-CZ	2.45	1.43	1.37
6	F	1	DTR	CZ2-CE2	2.38	1.46	1.41
6	F	4	PHA	CB-CG	2.07	1.56	1.51

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	F	5	A1D5E	O71-C72-N74	4.69	119.34	111.11
6	F	6	004	CB-CA-N	-4.49	101.65	112.40
6	F	5	A1D5E	CG-O71-C72	-3.81	111.02	116.48
6	F	5	A1D5E	O73-C72-N74	-3.01	120.35	124.96
6	F	5	A1D5E	O71-C72-O73	-2.80	120.31	124.53
6	F	1	DTR	CB-CA-C	-2.39	107.00	111.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	F	6	004	CG2-CB-CG1	2.26	121.11	118.29
6	F	3	TY5	C49-OH-CZ	-2.20	112.21	117.65

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	F	1	DTR	CA-CB-CG-CD1
6	F	3	TY5	C-CA-CB-CG
6	F	3	TY5	N-CA-CB-CG
6	F	5	A1D5E	O71-C72-N74-C75
6	F	5	A1D5E	O73-C72-N74-C75
6	F	5	A1D5E	N74-C72-O71-CG
6	F	5	A1D5E	O73-C72-O71-CG
6	F	3	TY5	CE2-CZ-OH-C49
6	F	3	TY5	CE1-CZ-OH-C49
6	F	3	TY5	CA-CB-CG-CD1
6	F	3	TY5	CA-CB-CG-CD2
6	F	5	A1D5E	C76-C75-N74-C72
6	F	6	004	C-CA-CB-CG2

There are no ring outliers.

4 monomers are involved in 38 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	F	3	TY5	18	0
6	F	6	004	16	0
6	F	1	DTR	2	0
6	F	4	PHA	3	0

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.

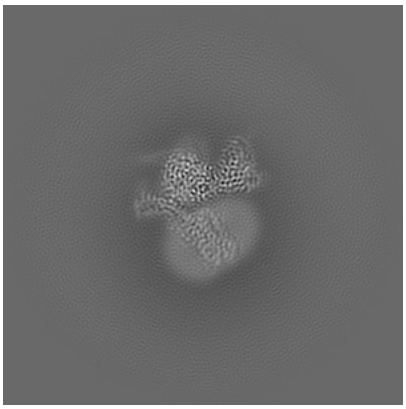
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-60650. These allow visual inspection of the internal detail of the map and identification of artifacts.

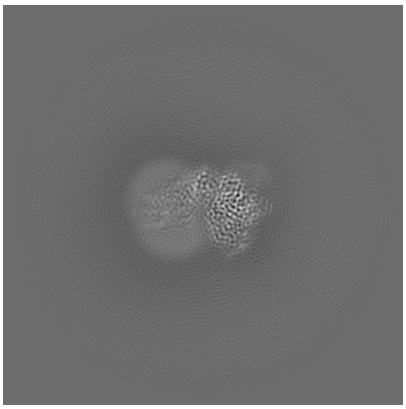
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

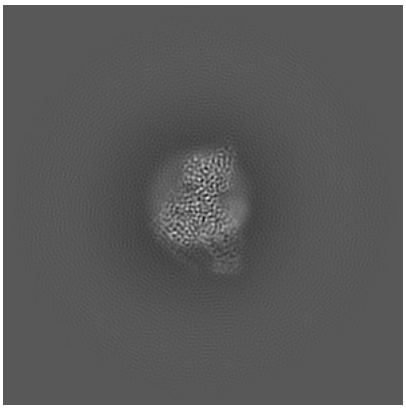
6.1.1 Primary map



X

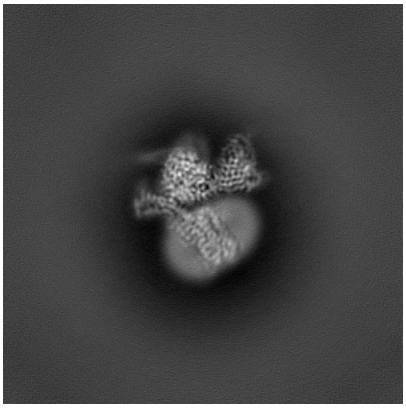


Y

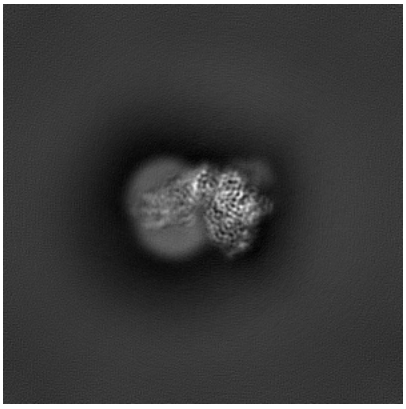


Z

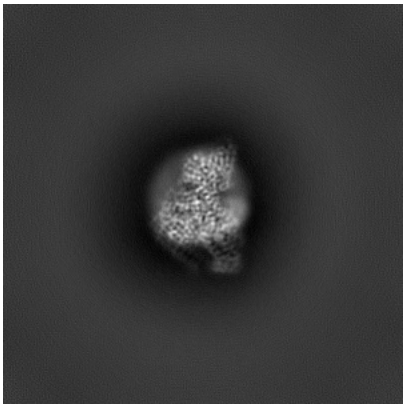
6.1.2 Raw map



X



Y

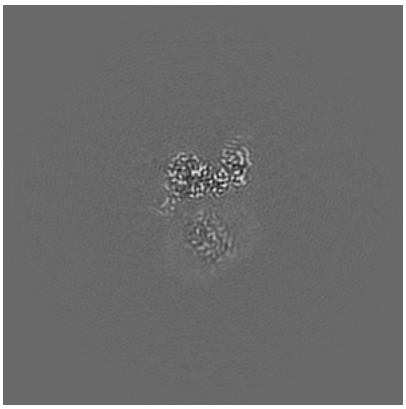


Z

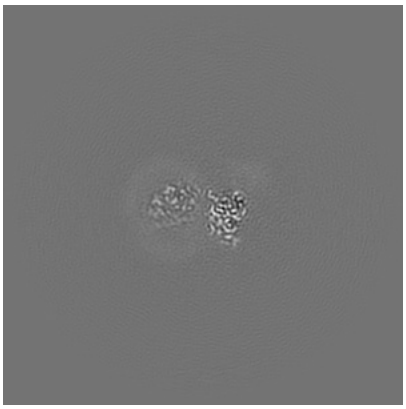
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

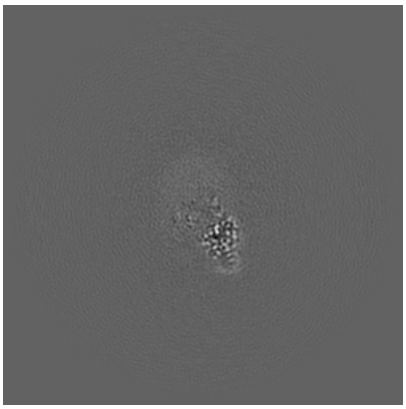
6.2.1 Primary map



X Index: 192

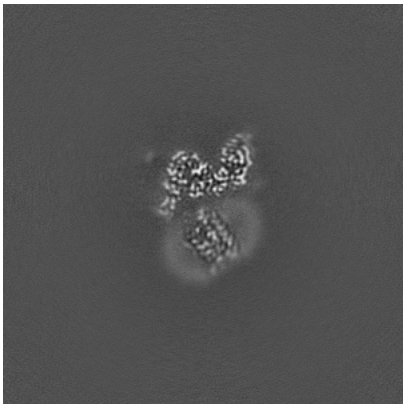


Y Index: 192

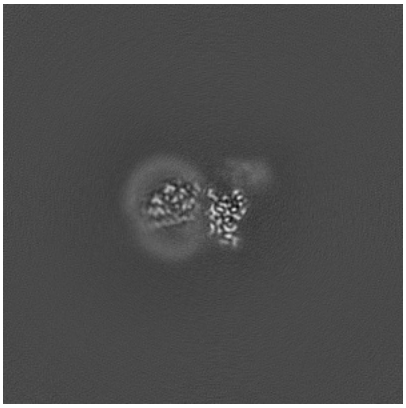


Z Index: 192

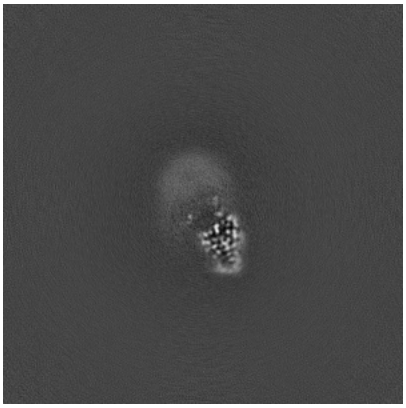
6.2.2 Raw map



X Index: 192



Y Index: 192

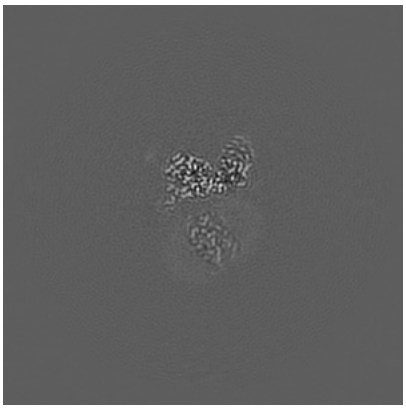


Z Index: 192

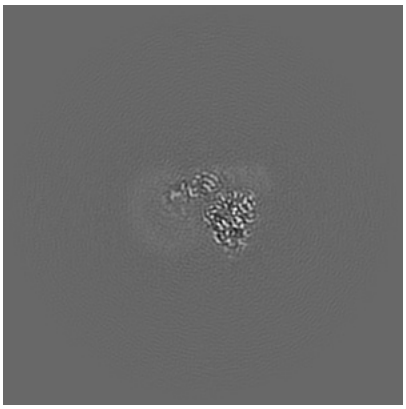
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

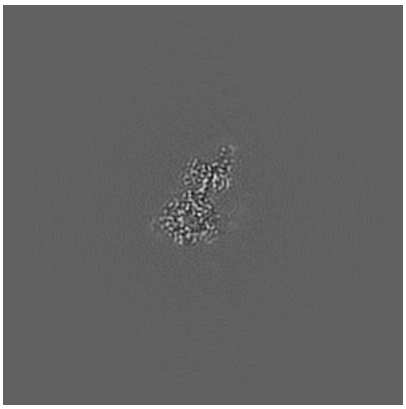
6.3.1 Primary map



X Index: 190

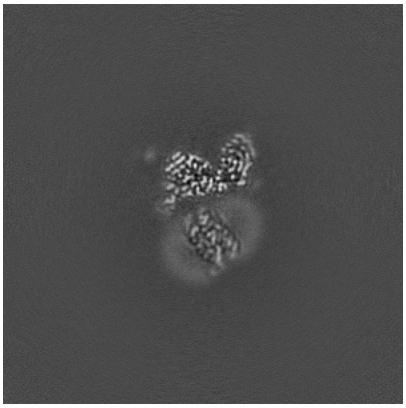


Y Index: 173

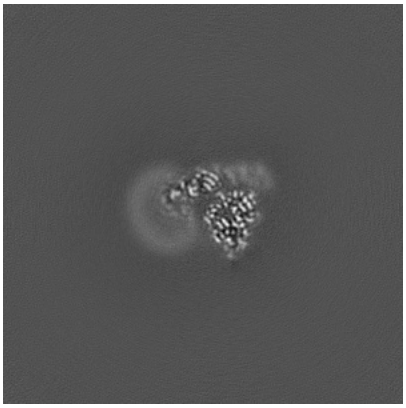


Z Index: 216

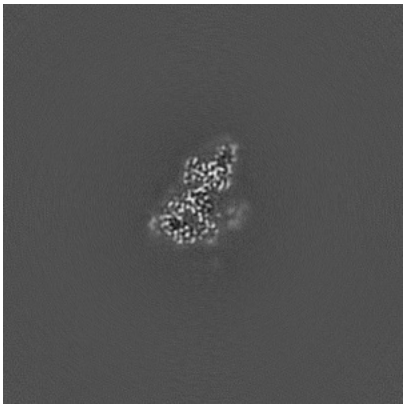
6.3.2 Raw map



X Index: 190



Y Index: 173

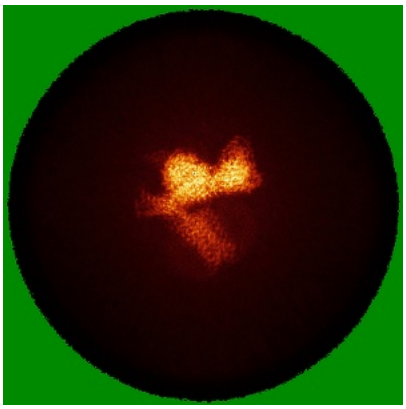


Z Index: 217

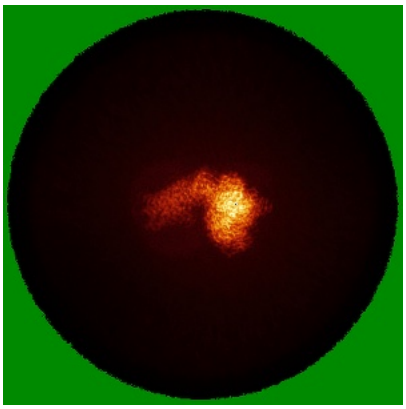
The images above show the largest variance slices of the map in three orthogonal directions.

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

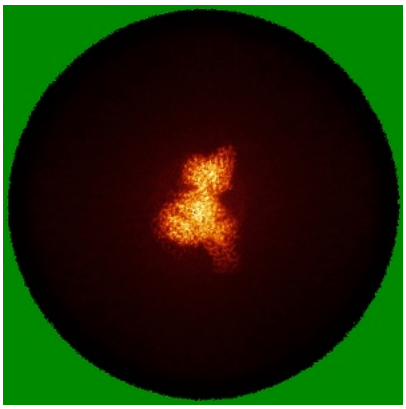
6.4.1 Primary map



X

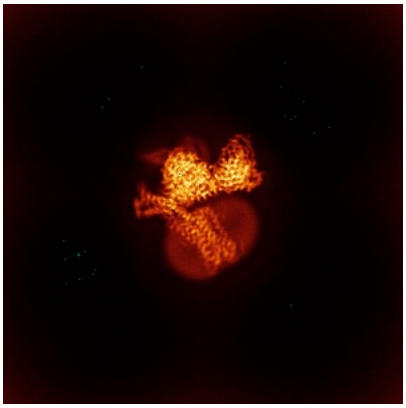


Y

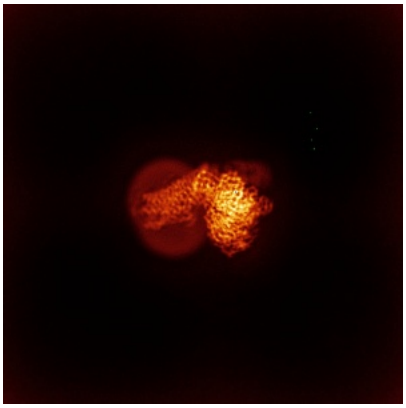


Z

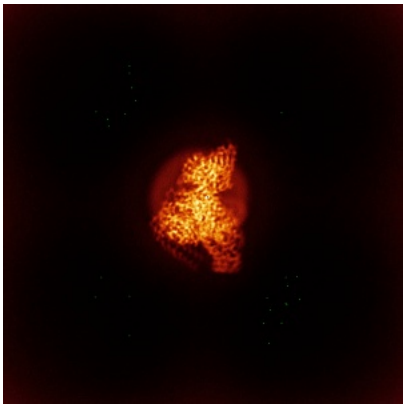
6.4.2 Raw map



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.

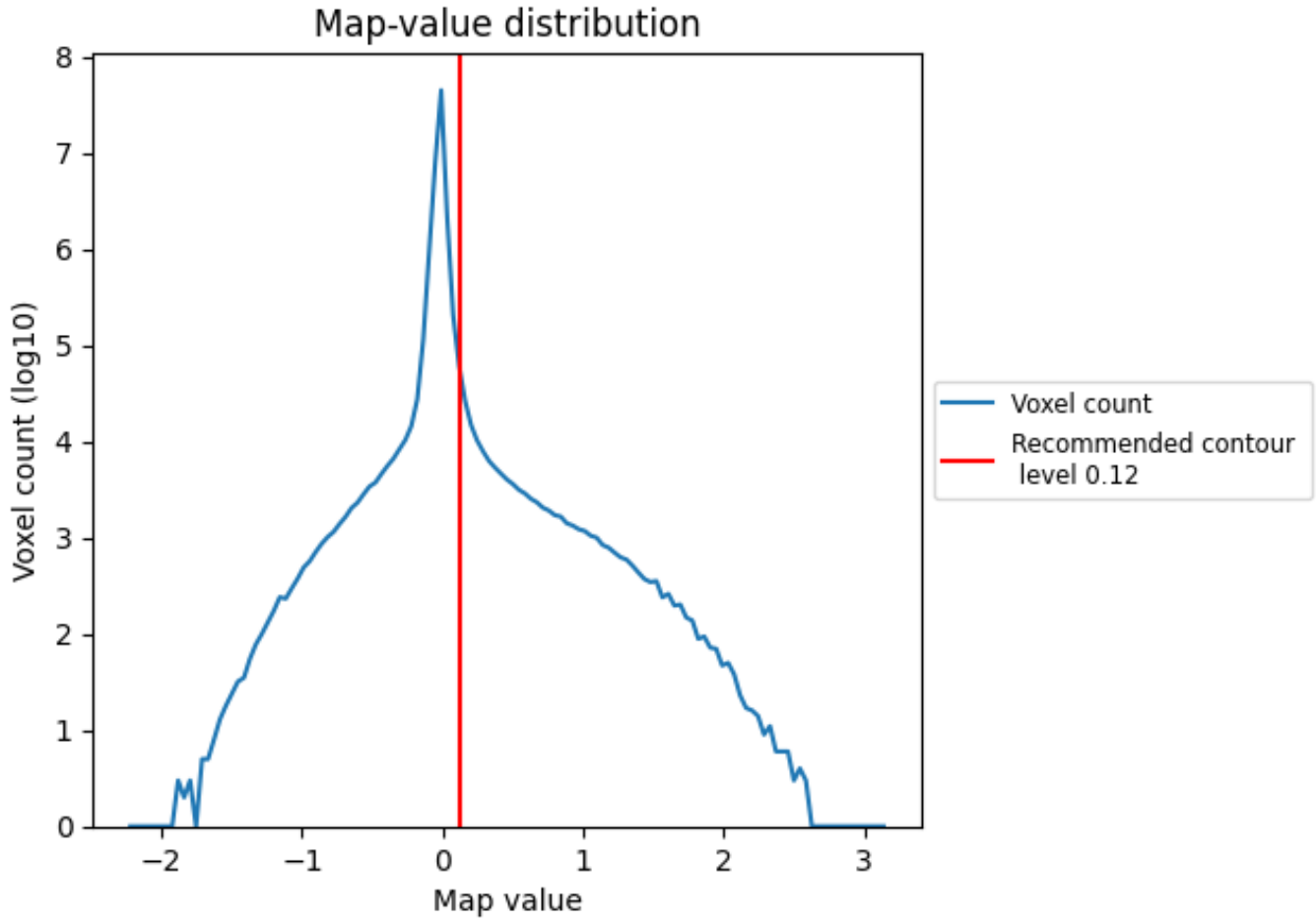
6.6 Mask visualisation (i)

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

7 Map analysis

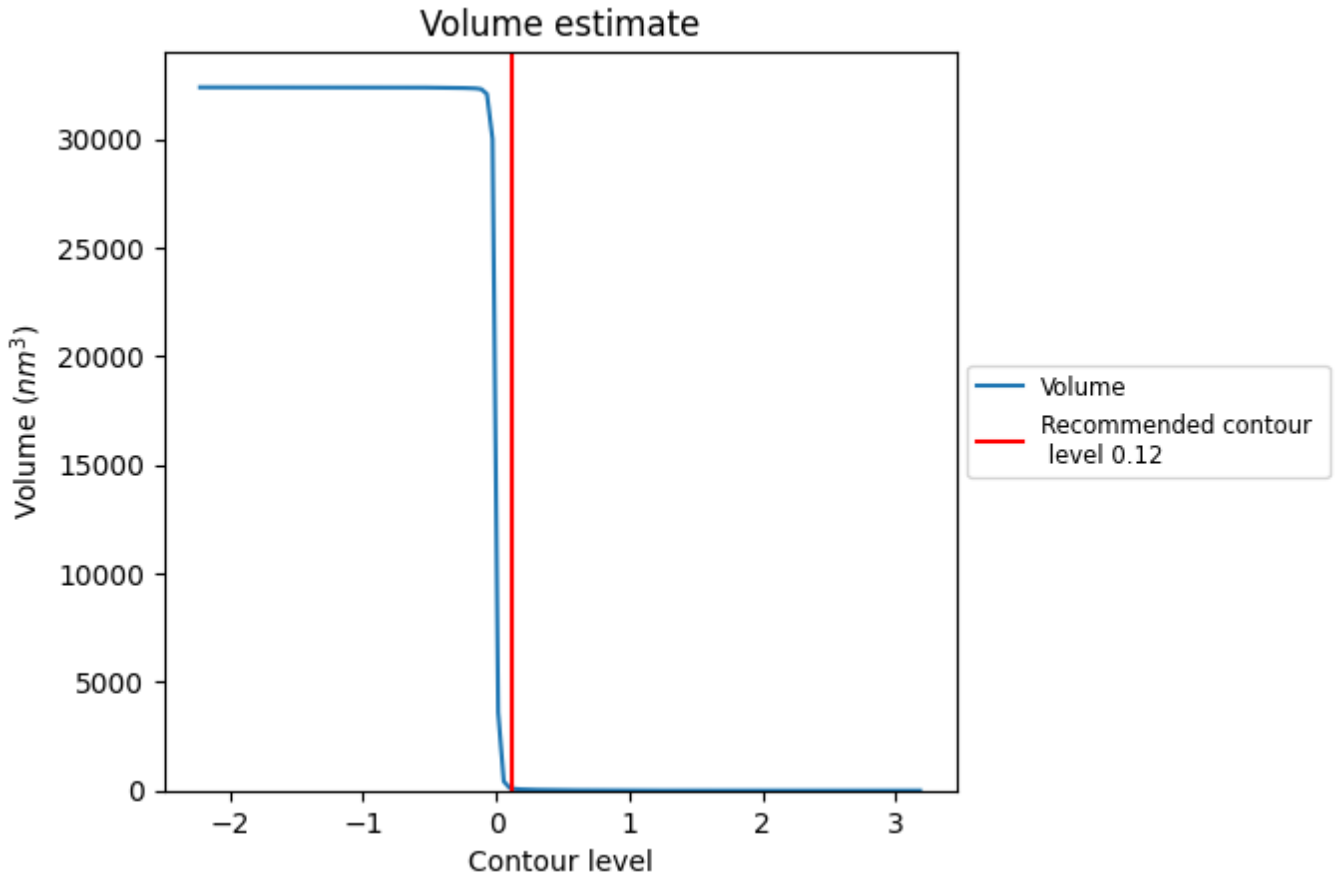
This section contains the results of statistical analysis of the map.

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.

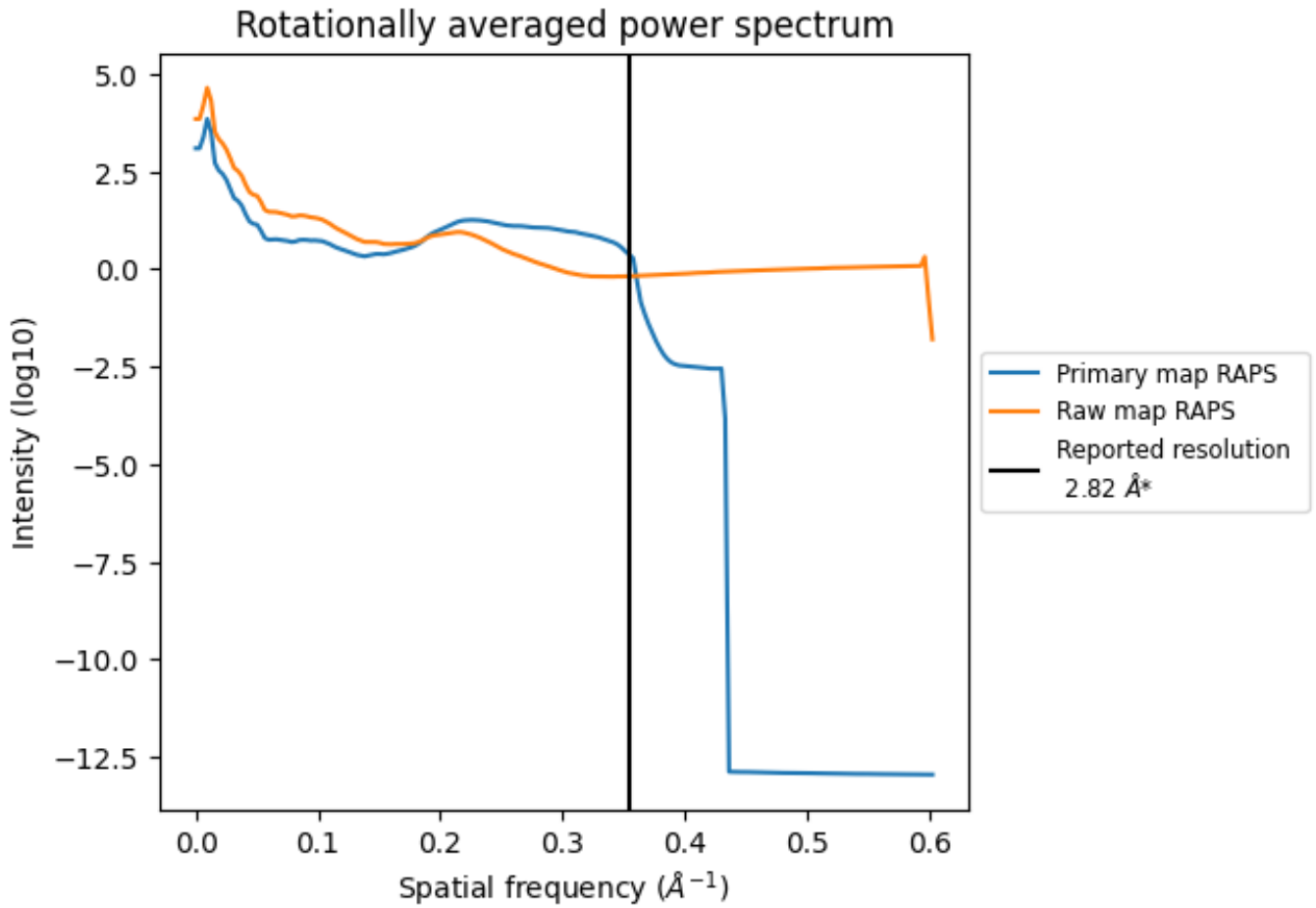
7.2 Volume estimate



The volume at the recommended contour level is 106 nm^3; this corresponds to an approximate mass of 95 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)

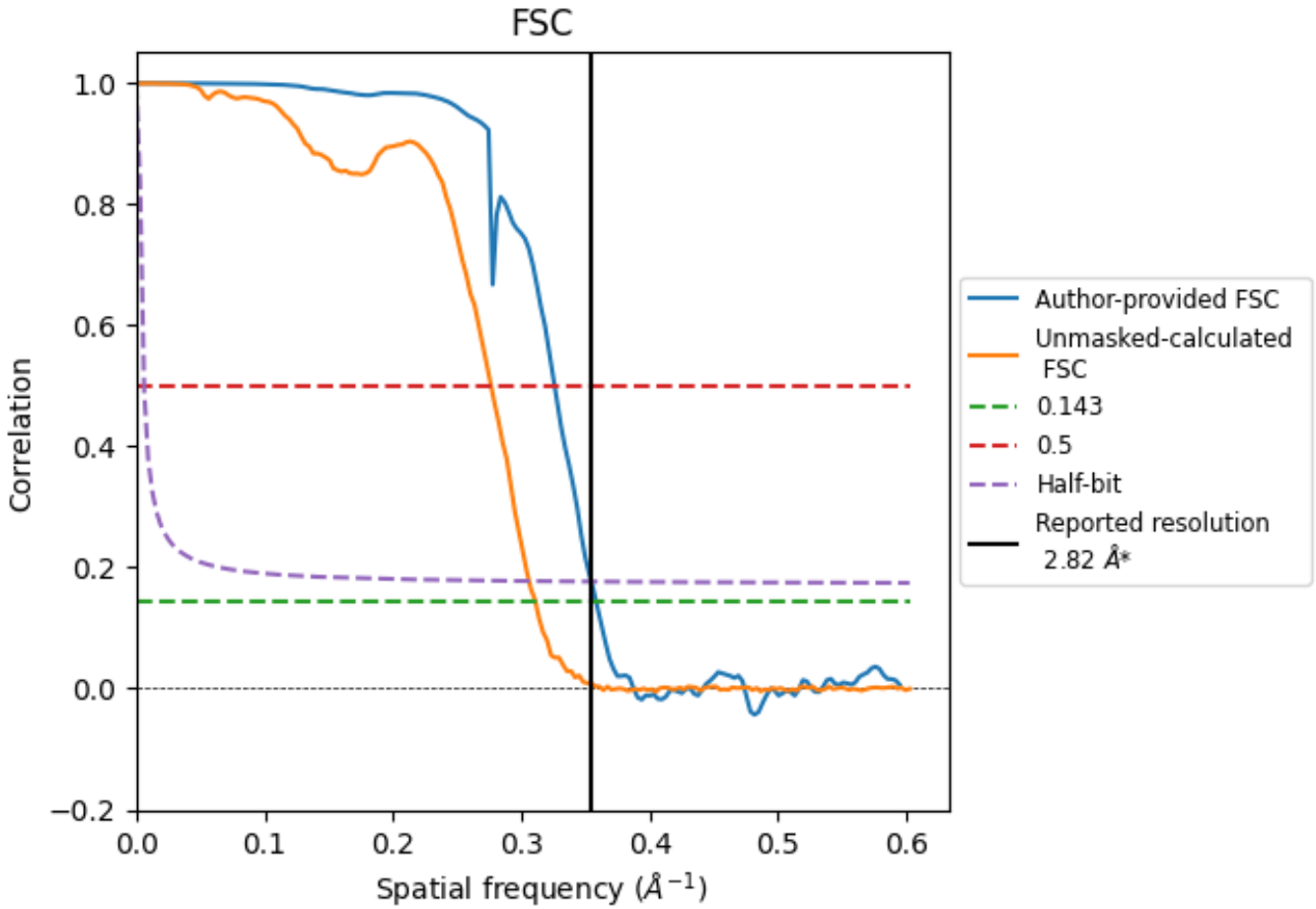


*Reported resolution corresponds to spatial frequency of 0.355 Å^{-1}

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 0.355 Å^{-1}

8.2 Resolution estimates (i)

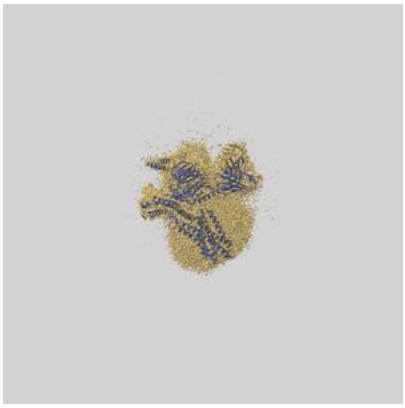
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.82	-	-
Author-provided FSC curve	2.79	3.07	2.82
Unmasked-calculated*	3.22	3.62	3.27

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.22 differs from the reported value 2.82 by more than 10 %

9 Map-model fit

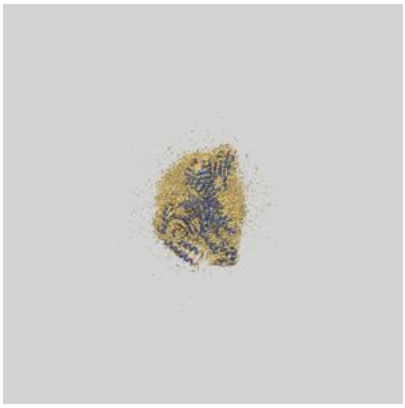
This section contains information regarding the fit between EMDB map EMD-60650 and PDB model 9IK8. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay



X

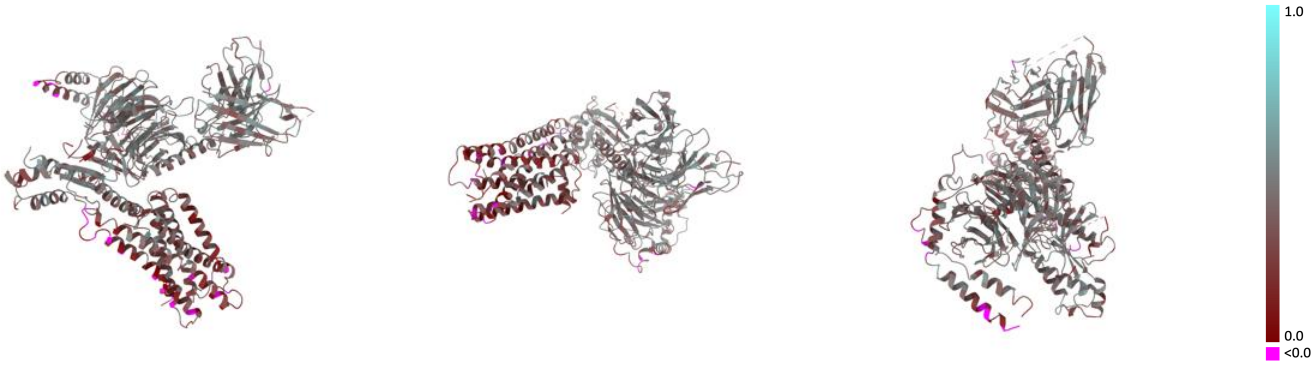
Y



Z

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

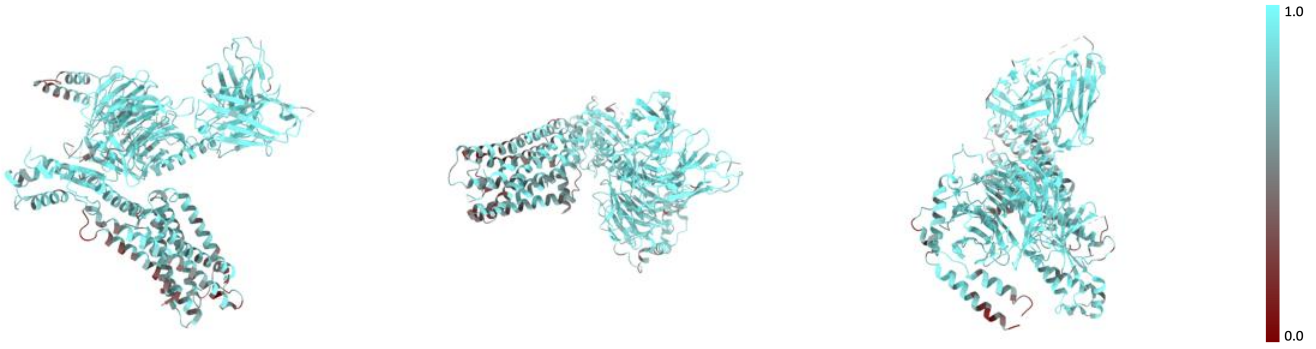
9.2 Q-score mapped to coordinate model (i)



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.

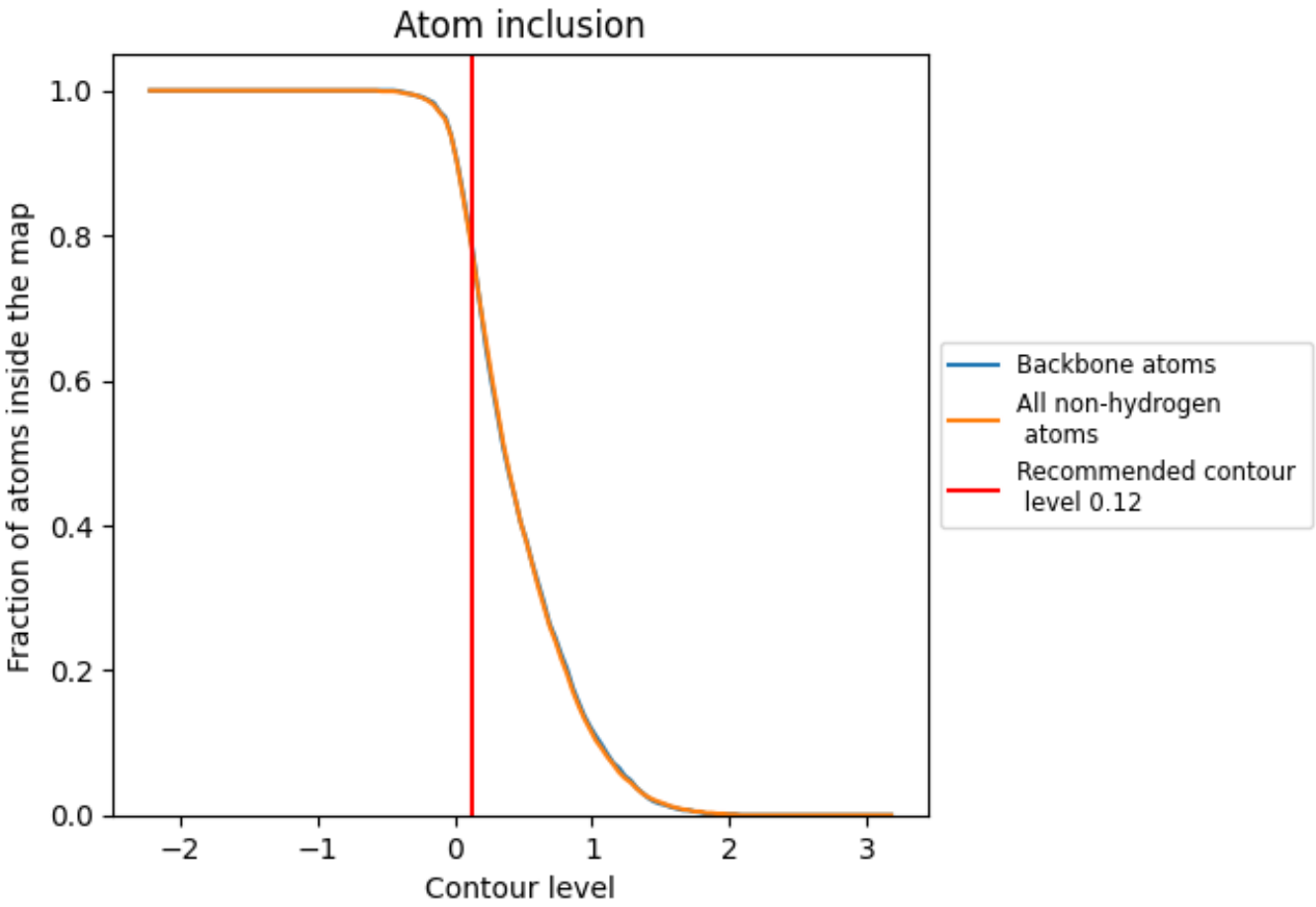
9.3 Atom inclusion mapped to coordinate model (i)



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, 79% of all backbone atoms, 79% 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 (0.12) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7860	0.3760
A	0.8200	0.4190
B	0.8500	0.4130
C	0.7530	0.3910
D	0.6360	0.2430
E	0.8740	0.4440
F	0.6930	0.3100