



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 08:06 PM UTC

PDB ID : 6B48 / pdb_00006b48
EMDB ID : EMD-7052
Title : Cryo-EM structure of Type I-F CRISPR crRNA-guided Csy surveillance complex with bound anti-CRISPR protein AcrF10
Authors : Guo, T.W.; Bartesaghi, A.; Yang, H.; Falconieri, V.; Rao, P.; Merk, A.; Fox, T.; Earl, L.; Patel, D.J.; Subramaniam, S.
Deposited on : 2017-09-25
Resolution : 3.60 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

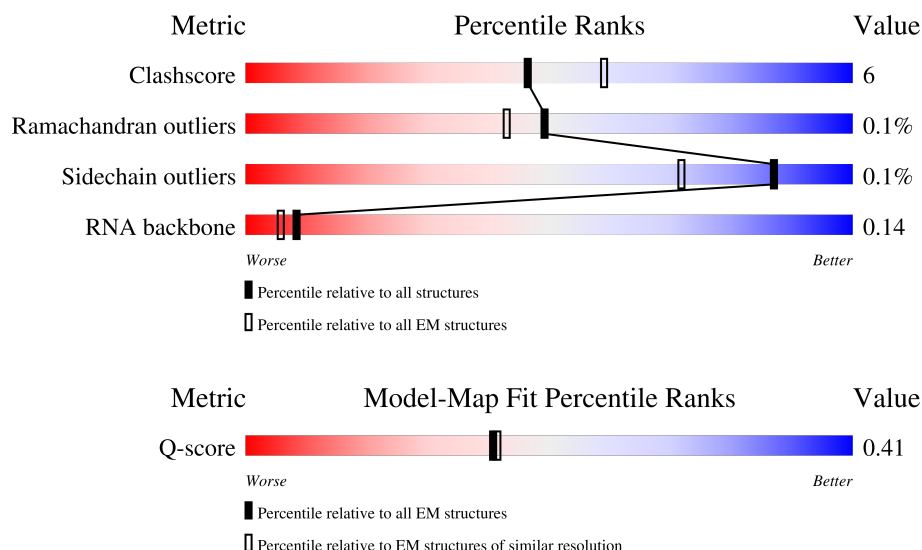
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13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

1 Overall quality at a glance

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

The reported resolution of this entry is 3.60 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	12797 (3.10 - 4.10)

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	436	55% 90% 7% .
2	B	329	22% 73% 20% 7%
3	C	344	85% 67% 18% 15%

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Mol	Chain	Length	Quality of chain
3	D	344	
3	E	344	
3	F	344	
3	G	344	
3	H	344	
4	K	99	
5	L	189	
6	M	60	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 22333 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 CRISPR-associated protein Csy1.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	424	Total	C	N	O	0	0
			2079	1145	473	461		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q02ML9
A	0	SER	-	expression tag	UNP Q02ML9

- Molecule 2 is a protein called CRISPR-associated protein Csy2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	305	Total	C	N	O	S	0	0
			2374	1504	440	425	5		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	MET	-	initiating methionine	UNP Q02MM0
B	0	ALA	-	expression tag	UNP Q02MM0

- Molecule 3 is a protein called CRISPR-associated protein Csy3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	293	Total	C	N	O	S	0	0
			2272	1430	409	431	2		
3	D	333	Total	C	N	O	S	0	0
			2554	1603	466	483	2		
3	E	334	Total	C	N	O	S	0	0
			2561	1611	466	482	2		
3	F	335	Total	C	N	O	S	0	0
			2566	1614	467	483	2		

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Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	333	Total	C	N	O	S	0	0
			2557	1608	466	481	2		
3	H	333	Total	C	N	O	S	0	0
			2554	1603	466	483	2		

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	MET	-	initiating methionine	UNP Q02MM1
C	0	ALA	-	expression tag	UNP Q02MM1
D	-1	MET	-	initiating methionine	UNP Q02MM1
D	0	ALA	-	expression tag	UNP Q02MM1
E	-1	MET	-	initiating methionine	UNP Q02MM1
E	0	ALA	-	expression tag	UNP Q02MM1
F	-1	MET	-	initiating methionine	UNP Q02MM1
F	0	ALA	-	expression tag	UNP Q02MM1
G	-1	MET	-	initiating methionine	UNP Q02MM1
G	0	ALA	-	expression tag	UNP Q02MM1
H	-1	MET	-	initiating methionine	UNP Q02MM1
H	0	ALA	-	expression tag	UNP Q02MM1

- Molecule 4 is a protein called Anti-CRISPR protein AcrF10.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	K	97	Total	C	N	O	S	Se	0	0
			786	496	119	165	1	5		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	0	GLY	-	expression tag	UNP A0A073KP86
K	1	SER	-	expression tag	UNP A0A073KP86
K	2	MSE	-	expression tag	UNP A0A073KP86

- Molecule 5 is a protein called CRISPR-associated endonuclease Cas6/Csy4.

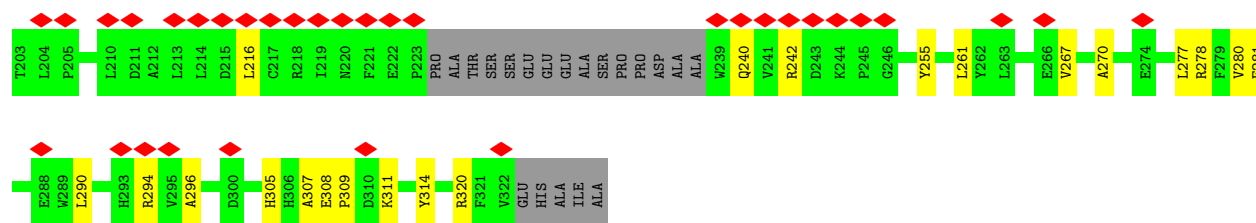
Mol	Chain	Residues	Atoms				AltConf	Trace
5	L	189	Total	C	N	O	0	0
			758	380	189	189		

There are 2 discrepancies between the modelled and reference sequences:

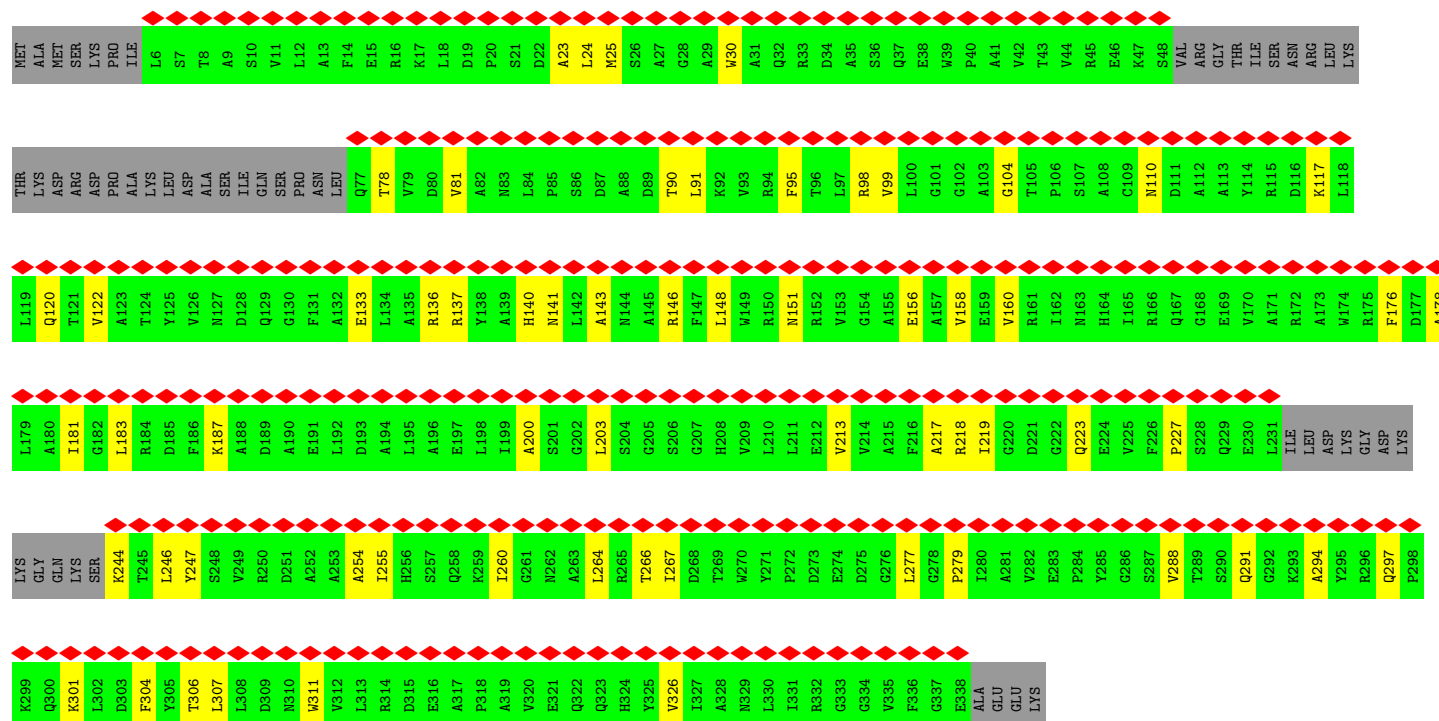
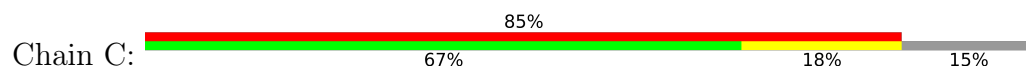
Chain	Residue	Modelled	Actual	Comment	Reference
L	-1	MET	-	initiating methionine	UNP Q02MM2
L	0	ALA	-	expression tag	UNP Q02MM2

- Molecule 6 is a RNA chain called *Pseudomonas aeruginosa* strain SMC4485 CRISPR repeat sequence.

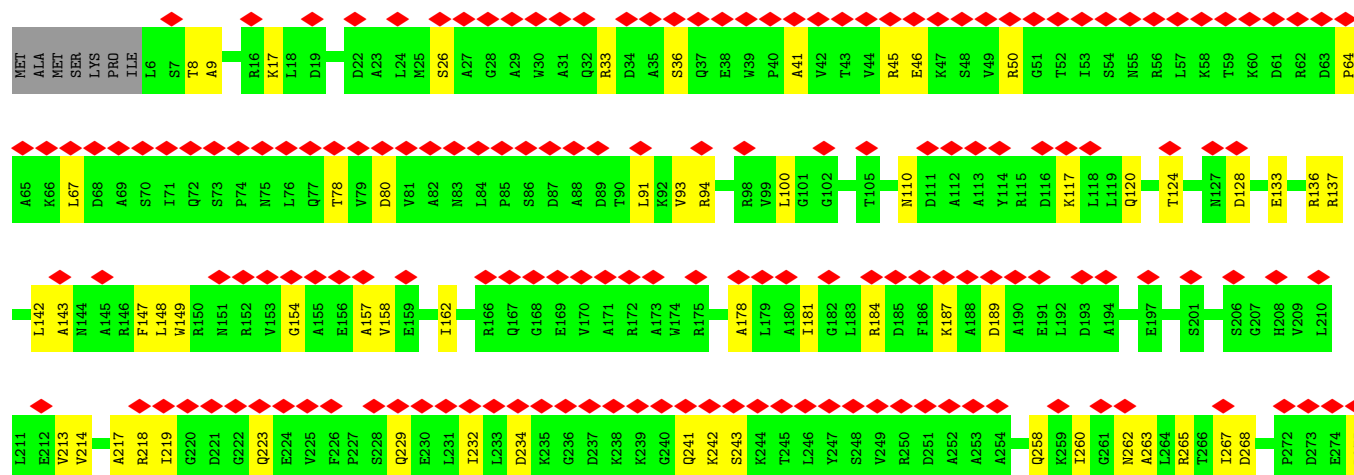
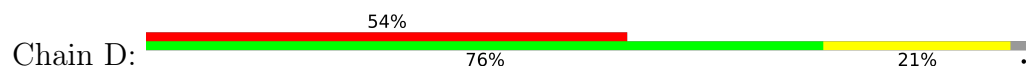
Mol	Chain	Residues	Atoms					AltConf	Trace
6	M	60	Total	C	N	O	P	0	0
			1272	569	223	421	59		

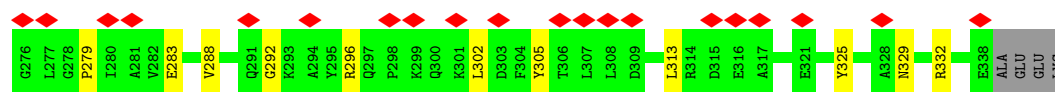


• Molecule 3: CRISPR-associated protein Csy3

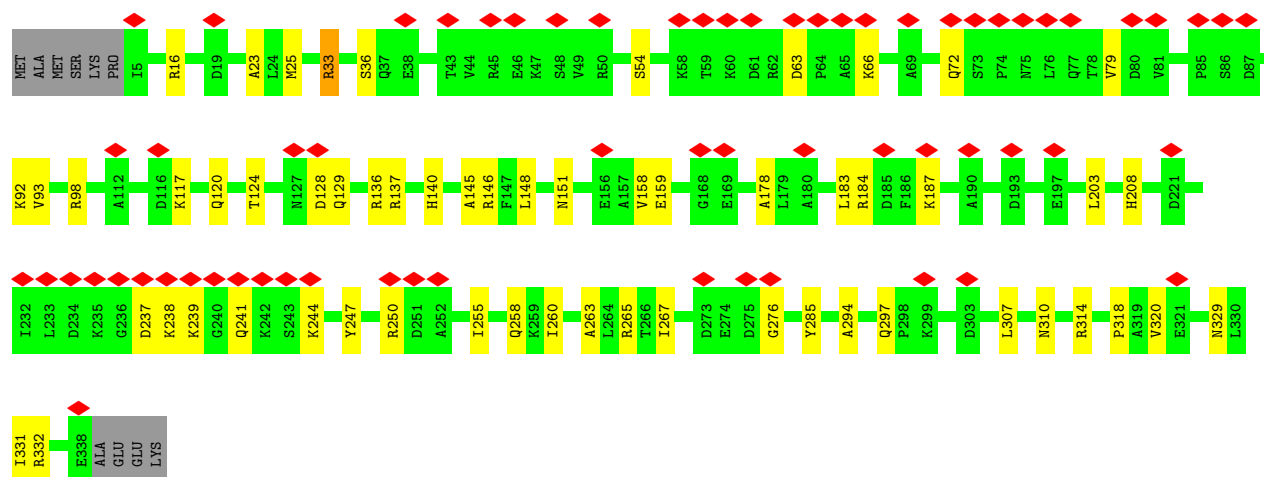
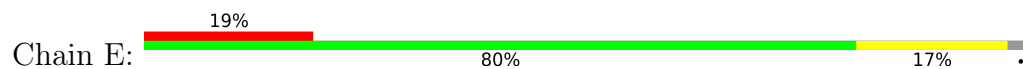


• Molecule 3: CRISPR-associated protein Csy3

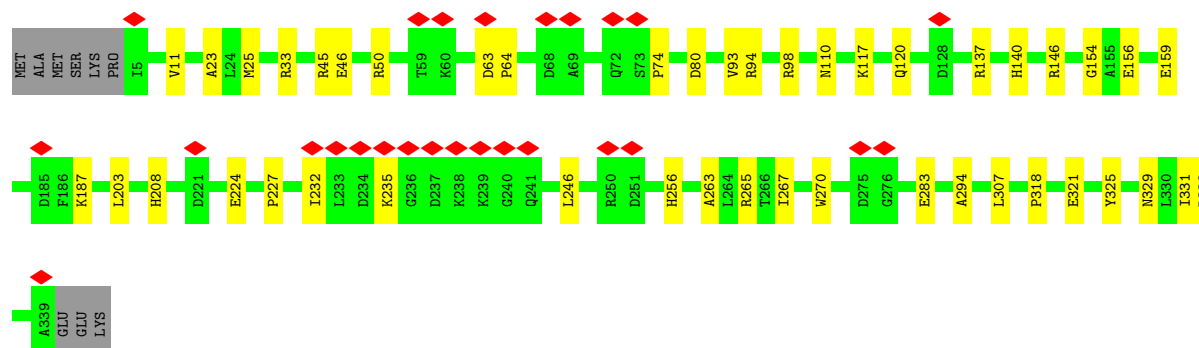
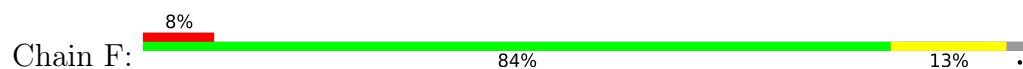




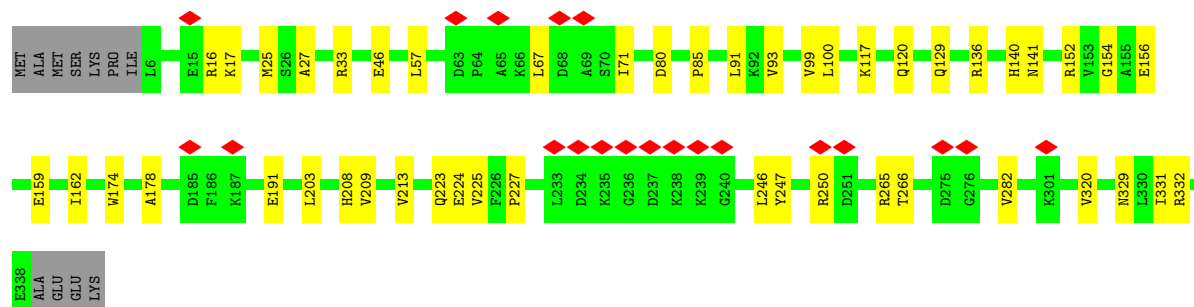
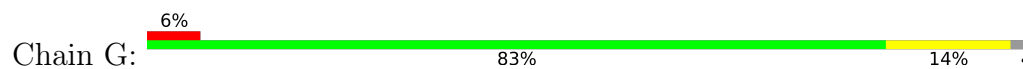
• Molecule 3: CRISPR-associated protein Csy3



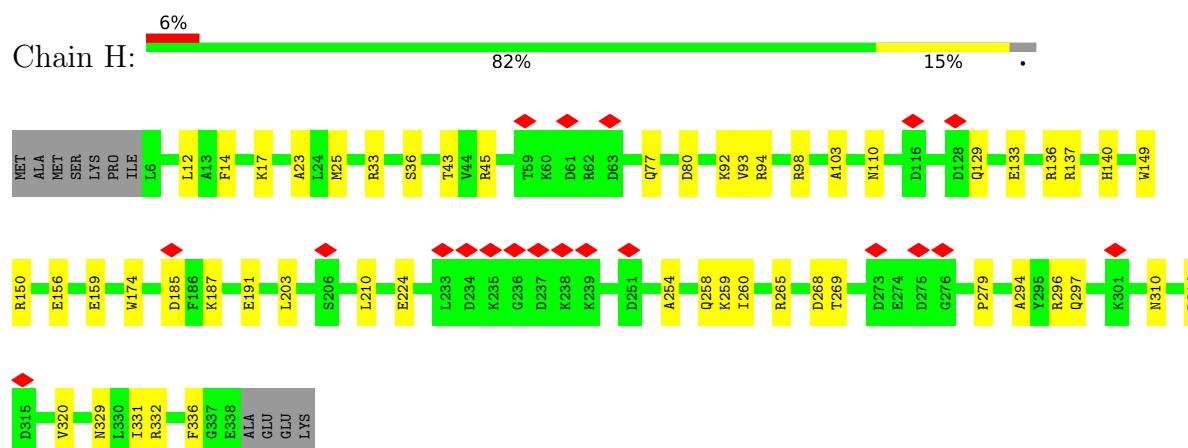
• Molecule 3: CRISPR-associated protein Csy3



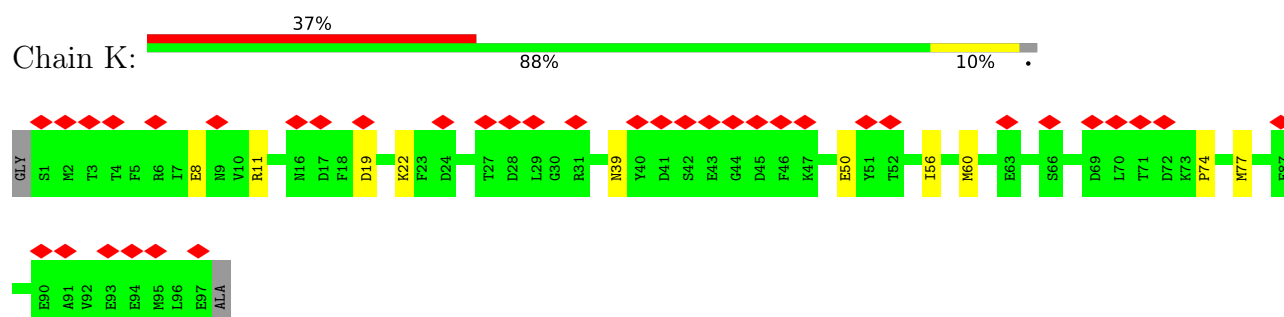
• Molecule 3: CRISPR-associated protein Csy3



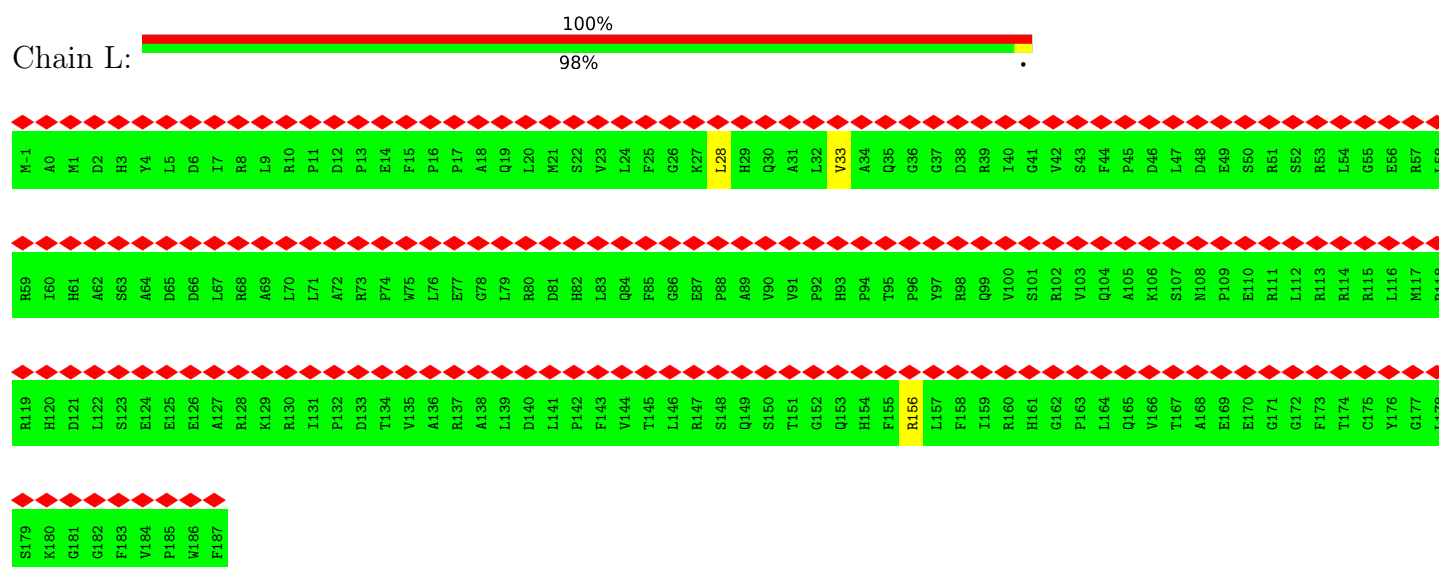
- Molecule 3: CRISPR-associated protein Csy3



- Molecule 4: Anti-CRISPR protein AcrF10



- Molecule 5: CRISPR-associated endonuclease Cas6/Csy4



- Molecule 6: Pseudomonas aeruginosa strain SMC4485 CRISPR repeat sequence





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	42498	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 ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	15.350	Depositor
Minimum map value	-9.633	Depositor
Average map value	0.018	Depositor
Map value standard deviation	0.748	Depositor
Recommended contour level	2.7	Depositor
Map size (Å)	199.92, 199.92, 199.92	wwPDB
Map dimensions	238, 238, 238	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.84, 0.84, 0.84	Depositor

5 Model quality

5.1 Standard geometry

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	0.24	0/2104	0.61	0/2715
2	B	0.31	0/2431	0.64	0/3310
3	C	0.17	0/2315	0.46	0/3143
3	D	0.25	0/2601	0.55	0/3532
3	E	0.31	0/2608	0.58	0/3540
3	F	0.35	0/2613	0.58	0/3547
3	G	0.36	0/2604	0.61	0/3533
3	H	0.34	0/2601	0.62	0/3532
4	K	0.31	0/793	0.68	2/1066 (0.2%)
5	L	0.11	0/757	0.33	0/946
6	M	0.28	0/1420	0.43	0/2212
All	All	0.29	0/22847	0.57	2/31076 (0.0%)

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	1
2	B	0	2
3	E	0	1
All	All	0	4

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	K	8	GLU	CA-C-N	5.06	131.20	121.54
4	K	8	GLU	C-N-CA	5.06	131.20	121.54

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	222	ARG	Peptide
2	B	105	LEU	Peptide
2	B	280	VAL	Peptide
3	E	285	TYR	Peptide

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	A	2079	0	1094	21	0
2	B	2374	0	2345	43	0
3	C	2272	0	2232	39	0
3	D	2554	0	2522	45	0
3	E	2561	0	2542	37	0
3	F	2566	0	2547	28	0
3	G	2557	0	2542	32	0
3	H	2554	0	2522	34	0
4	K	786	0	749	5	0
5	L	758	0	209	2	0
6	M	1272	0	644	19	0
All	All	22333	0	19948	259	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 6.

All (259) 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:G:25:MET:HG2	3:G:93:VAL:HG22	1.78	0.65
2:B:148:CYS:SG	2:B:149:ASN:N	2.70	0.64
3:D:262:ASN:ND2	6:M:32:G:OP2	2.30	0.64
3:F:232:ILE:HG13	3:F:235:LYS:HB2	1.80	0.63
2:B:25:SER:HB3	2:B:28:THR:HG22	1.82	0.60
3:G:203:LEU:HD22	3:G:331:ILE:HD11	1.82	0.60
3:H:33:ARG:NH1	3:H:159:GLU:OE1	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:33:ARG:NH1	3:E:159:GLU:OE1	2.35	0.60
2:B:270:ALA:O	3:H:110:ASN:ND2	2.35	0.59
3:G:129:GLN:HE22	3:G:320:VAL:HG13	1.67	0.59
1:A:92:PRO:O	2:B:194:GLN:NE2	2.36	0.59
2:B:18:GLN:HE21	2:B:143:SER:HB3	1.68	0.58
1:A:176:LYS:NZ	6:M:4:A:OP2	2.35	0.58
3:F:33:ARG:NH1	3:F:159:GLU:OE1	2.36	0.58
3:C:78:THR:O	3:C:244:LYS:N	2.36	0.58
3:G:80:ASP:OD1	3:G:80:ASP:N	2.35	0.58
3:D:133:GLU:OE1	3:D:136:ARG:NH2	2.36	0.58
3:D:9:ALA:HB1	3:D:110:ASN:HB2	1.84	0.58
3:C:247:TYR:OH	3:D:45:ARG:NH1	2.36	0.57
3:D:26:SER:HA	3:D:41:ALA:HA	1.86	0.57
3:C:146:ARG:HH21	3:C:183:LEU:HG	1.69	0.57
3:C:136:ARG:O	3:C:140:HIS:ND1	2.38	0.57
3:C:133:GLU:OE1	3:C:136:ARG:NH2	2.38	0.57
3:D:124:THR:O	3:D:128:ASP:HB2	2.04	0.57
5:L:156:ARG:O	6:M:46:C:N4	2.37	0.57
3:G:33:ARG:NH1	3:G:159:GLU:OE1	2.39	0.56
3:E:310:ASN:OD1	3:E:314:ARG:NH1	2.39	0.56
3:C:98:ARG:NH1	3:D:154:GLY:O	2.38	0.56
2:B:261:LEU:HD12	2:B:278:ARG:HD2	1.87	0.56
3:D:91:LEU:HB3	3:D:217:ALA:HB3	1.88	0.56
3:F:98:ARG:NH1	3:G:154:GLY:O	2.39	0.56
2:B:255:TYR:HB2	2:B:314:TYR:HB2	1.87	0.56
3:G:17:LYS:HG3	3:G:100:LEU:HB2	1.87	0.56
2:B:67:GLU:HG2	2:B:104:HIS:HB2	1.88	0.56
3:G:152:ARG:NH2	3:G:178:ALA:O	2.37	0.56
3:H:129:GLN:HE22	3:H:320:VAL:HG13	1.71	0.55
3:E:98:ARG:NH1	3:F:154:GLY:O	2.40	0.55
2:B:88:LEU:HB3	2:B:92:GLY:HA2	1.88	0.55
3:D:33:ARG:HA	3:D:36:SER:HB3	1.87	0.55
2:B:83:LEU:HD11	2:B:99:GLU:HG2	1.88	0.55
3:C:81:VAL:HG22	3:C:246:LEU:HD11	1.88	0.55
3:C:143:ALA:HB1	3:C:181:ILE:HG21	1.89	0.55
3:H:310:ASN:OD1	3:H:314:ARG:NH1	2.40	0.55
3:E:23:ALA:HB2	3:E:260:ILE:HD11	1.88	0.55
3:C:25:MET:HB3	3:C:91:LEU:HD11	1.87	0.55
3:H:133:GLU:OE1	3:H:136:ARG:NH2	2.41	0.54
3:E:237:ASP:OD2	3:E:241:GLN:NE2	2.41	0.54
3:E:137:ARG:HB3	3:E:267:ILE:HD12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:41:VAL:HG21	2:B:59:VAL:HB	1.88	0.54
2:B:148:CYS:SG	3:H:94:ARG:NH2	2.81	0.54
3:G:99:VAL:HB	3:G:209:VAL:HG13	1.89	0.54
3:E:265:ARG:NH2	6:M:26:G:OP2	2.41	0.54
3:D:283:GLU:O	3:D:325:TYR:OH	2.26	0.53
3:G:208:HIS:ND1	3:H:156:GLU:OE2	2.38	0.53
1:A:195:PRO:HB3	2:B:281:GLU:HG2	1.89	0.53
3:C:117:LYS:HD2	3:C:120:GLN:HE21	1.74	0.53
3:G:162:ILE:HG12	3:G:213:VAL:HG22	1.90	0.53
3:E:140:HIS:NE2	3:E:187:LYS:O	2.38	0.53
3:G:16:ARG:NH2	3:H:224:GLU:OE1	2.41	0.53
3:G:27:ALA:HB3	3:G:85:PRO:HD2	1.91	0.53
1:A:251:ILE:HB	1:A:254:LEU:HD12	1.90	0.53
2:B:308:GLU:HG3	2:B:311:LYS:HB2	1.91	0.53
3:E:247:TYR:OH	3:F:45:ARG:NH1	2.38	0.53
3:H:258:GLN:NE2	6:M:9:U:OP1	2.41	0.53
3:H:150:ARG:NH1	6:M:11:C:OP1	2.38	0.53
1:A:390:ARG:O	3:E:238:LYS:NZ	2.42	0.53
2:B:124:GLN:HE22	2:B:128:ARG:HD3	1.73	0.53
2:B:146:PRO:HG3	3:H:210:LEU:HD11	1.91	0.53
3:E:36:SER:O	3:E:92:LYS:NZ	2.42	0.53
3:G:16:ARG:HB2	6:M:12:A:H5"	1.90	0.53
3:E:208:HIS:ND1	3:F:156:GLU:OE2	2.41	0.53
3:C:227:PRO:HG2	3:C:246:LEU:HD22	1.91	0.52
3:E:117:LYS:HD2	3:E:120:GLN:HE21	1.73	0.52
5:L:28:LEU:O	5:L:33:VAL:N	2.43	0.52
3:G:227:PRO:HG2	3:G:246:LEU:HD22	1.91	0.52
1:A:224:ARG:NH2	1:A:226:GLU:OE2	2.38	0.52
3:D:8:THR:HG23	3:D:313:LEU:HD11	1.91	0.52
2:B:5:ASP:HA	2:B:320:ARG:HH12	1.73	0.52
3:D:162:ILE:HG12	3:D:213:VAL:HG22	1.92	0.52
3:E:184:ARG:NH2	3:E:276:GLY:O	2.40	0.52
3:E:258:GLN:NE2	6:M:27:U:OP1	2.43	0.52
2:B:10:LEU:HB3	2:B:111:LEU:HB2	1.92	0.51
3:E:294:ALA:O	3:E:297:GLN:NE2	2.37	0.51
3:H:14:PHE:O	6:M:5:G:O2'	2.22	0.51
3:H:80:ASP:OD1	3:H:80:ASP:N	2.39	0.51
3:C:141:ASN:ND2	3:C:264:LEU:O	2.43	0.51
2:B:62:VAL:HG11	2:B:180:LEU:HD13	1.92	0.51
3:F:80:ASP:OD1	3:F:80:ASP:N	2.38	0.51
3:D:117:LYS:HD3	3:D:120:GLN:HE21	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:250:ARG:NH2	3:H:43:THR:OG1	2.43	0.51
3:E:36:SER:HB2	3:E:92:LYS:HD2	1.93	0.51
4:K:39:ASN:ND2	4:K:50:GLU:OE1	2.41	0.51
3:F:25:MET:HG2	3:F:93:VAL:HG22	1.92	0.51
3:D:46:GLU:HG2	3:D:80:ASP:HB3	1.91	0.51
3:E:148:LEU:HD22	3:E:151:ASN:HB2	1.92	0.51
3:H:23:ALA:HB2	3:H:260:ILE:HD11	1.93	0.51
3:H:25:MET:HG2	3:H:93:VAL:HG22	1.91	0.51
1:A:121:LEU:O	1:A:131:LEU:N	2.43	0.50
1:A:181:PRO:HG2	2:B:305:HIS:CE1	2.47	0.50
3:D:94:ARG:HG3	3:D:214:VAL:HG22	1.93	0.50
3:F:208:HIS:ND1	3:G:156:GLU:OE2	2.43	0.50
1:A:254:LEU:HD22	1:A:260:GLY:HA2	1.94	0.50
3:C:23:ALA:HA	3:C:95:PHE:HB3	1.93	0.50
2:B:267:VAL:HG11	2:B:277:LEU:HD22	1.94	0.50
3:F:256:HIS:ND1	6:M:21:U:OP1	2.41	0.50
2:B:21:ASN:ND2	6:M:4:A:OP1	2.41	0.50
3:C:90:THR:HG22	3:C:218:ARG:HG2	1.93	0.49
3:C:266:THR:HG22	3:C:279:PRO:HB2	1.93	0.49
3:E:203:LEU:HD22	3:E:331:ILE:HD11	1.93	0.49
3:C:304:PHE:HE1	3:C:326:VAL:HG13	1.78	0.49
3:D:50:ARG:O	6:M:37:C:O2'	2.31	0.49
3:H:203:LEU:HD22	3:H:331:ILE:HD11	1.94	0.49
3:E:146:ARG:HE	3:E:183:LEU:HD11	1.76	0.49
3:F:137:ARG:HB3	3:F:267:ILE:HD12	1.94	0.49
3:H:17:LYS:HE3	3:H:103:ALA:HA	1.93	0.49
2:B:173:GLN:NE2	2:B:296:ALA:O	2.46	0.49
1:A:46:THR:O	1:A:52:TRP:N	2.46	0.49
1:A:174:LEU:HA	6:M:5:G:H1	1.76	0.49
3:H:133:GLU:OE2	3:H:137:ARG:NE	2.46	0.49
3:E:136:ARG:O	3:E:140:HIS:ND1	2.46	0.49
2:B:196:HIS:HA	2:B:199:THR:HG22	1.94	0.49
3:F:227:PRO:HG2	3:F:246:LEU:HD22	1.94	0.49
3:C:91:LEU:HB3	3:C:217:ALA:HB3	1.96	0.48
3:D:265:ARG:CZ	3:D:332:ARG:HE	2.26	0.48
3:D:288:VAL:O	3:D:292:GLY:N	2.46	0.48
3:D:219:ILE:HG22	3:D:223:GLN:HB3	1.96	0.48
3:F:146:ARG:HG3	3:F:263:ALA:HB1	1.95	0.48
3:F:117:LYS:HD2	3:F:120:GLN:HE21	1.78	0.48
4:K:19:ASP:N	4:K:19:ASP:OD1	2.47	0.48
4:K:56:ILE:HG22	4:K:60:MSE:HG3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:110:ASN:ND2	3:E:72:GLN:OE1	2.46	0.48
3:E:145:ALA:HB1	3:E:148:LEU:HD12	1.96	0.48
3:F:46:GLU:HG2	3:F:80:ASP:HB3	1.95	0.48
3:G:247:TYR:OH	3:H:45:ARG:NH1	2.39	0.48
3:C:294:ALA:O	3:C:297:GLN:NE2	2.46	0.48
3:D:229:GLN:NE2	3:D:243:SER:O	2.47	0.48
3:C:227:PRO:HA	3:C:255:ILE:HG12	1.95	0.47
2:B:17:ILE:HG22	2:B:144:ILE:HG12	1.96	0.47
3:C:30:TRP:HB2	3:C:218:ARG:HH12	1.78	0.47
3:F:283:GLU:O	3:F:325:TYR:OH	2.33	0.47
1:A:179:TYR:HA	1:A:189:LEU:HA	1.97	0.47
1:A:266:PRO:O	2:B:28:THR:OG1	2.27	0.47
3:F:203:LEU:HD22	3:F:331:ILE:HD11	1.97	0.47
3:F:329:ASN:HA	3:F:332:ARG:HG2	1.96	0.47
3:D:46:GLU:OE1	3:D:241:GLN:NE2	2.48	0.47
3:D:158:VAL:HB	3:D:178:ALA:HB3	1.97	0.47
3:E:23:ALA:HB3	3:E:255:ILE:HB	1.96	0.47
3:F:140:HIS:NE2	3:F:187:LYS:O	2.38	0.47
3:C:104:GLY:HA2	3:C:122:VAL:HG11	1.96	0.47
3:D:93:VAL:HG21	3:D:148:LEU:HD21	1.96	0.46
3:D:329:ASN:HA	3:D:332:ARG:HG2	1.96	0.46
2:B:17:ILE:HD13	2:B:139:LEU:HD12	1.97	0.46
3:E:158:VAL:HB	3:E:178:ALA:HB3	1.98	0.46
3:D:296:ARG:HG2	3:D:302:LEU:HB3	1.97	0.46
4:K:11:ARG:HB2	4:K:22:LYS:HB2	1.98	0.46
3:C:23:ALA:HB2	3:C:260:ILE:HD11	1.97	0.46
3:C:307:LEU:O	3:C:311:TRP:HB2	2.15	0.46
3:H:329:ASN:HA	3:H:332:ARG:HG2	1.97	0.46
3:H:77:GLN:NE2	6:M:13:C:N3	2.63	0.46
2:B:135:GLY:O	3:H:98:ARG:NH1	2.48	0.46
3:C:148:LEU:HB3	3:C:151:ASN:HB2	1.97	0.46
3:D:265:ARG:HH22	6:M:31:C:H4'	1.81	0.46
3:G:46:GLU:HG2	3:G:80:ASP:HB3	1.97	0.46
3:D:64:PRO:HA	3:D:67:LEU:HD13	1.97	0.45
3:D:143:ALA:HB1	3:D:181:ILE:HD13	1.98	0.45
3:G:117:LYS:HD2	3:G:120:GLN:HE21	1.81	0.45
3:E:329:ASN:HA	3:E:332:ARG:HG2	1.99	0.45
3:C:99:VAL:HG11	3:C:203:LEU:HD23	1.99	0.45
1:A:178:LEU:O	1:A:190:LEU:N	2.48	0.45
1:A:180:PHE:HE2	2:B:307:ALA:HB2	1.81	0.45
1:A:202:VAL:HG11	2:B:27:LEU:HD22	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:91:LEU:HD13	3:G:225:VAL:HG11	1.99	0.45
3:D:187:LYS:NZ	3:D:189:ASP:OD2	2.35	0.45
2:B:290:LEU:HD22	2:B:294:ARG:HG3	1.98	0.45
3:C:160:VAL:HB	3:C:176:PHE:HB2	1.98	0.45
3:D:137:ARG:NH1	3:D:267:ILE:O	2.50	0.45
3:H:294:ALA:O	3:H:297:GLN:NE2	2.38	0.45
3:C:158:VAL:HB	3:C:178:ALA:HB3	2.00	0.44
3:C:307:LEU:O	3:C:311:TRP:CB	2.65	0.44
3:E:25:MET:HG2	3:E:93:VAL:HG22	1.97	0.44
3:H:265:ARG:NH1	3:H:332:ARG:HE	2.15	0.44
3:H:36:SER:O	3:H:92:LYS:NZ	2.36	0.44
3:E:79:VAL:HG13	3:E:244:LYS:HB2	1.99	0.44
1:A:240:ILE:HD13	1:A:263:TRP:HB3	2.00	0.44
3:E:129:GLN:HE22	3:E:320:VAL:HG13	1.81	0.44
3:F:270:TRP:CD1	3:F:321:GLU:HB2	2.52	0.44
1:A:210:ARG:NH2	1:A:235:TYR:H	2.15	0.44
3:H:174:TRP:HD1	3:H:191:GLU:HG3	1.83	0.44
3:C:219:ILE:HG22	3:C:223:GLN:HB3	1.99	0.44
4:K:74:PRO:HD2	4:K:77:MSE:HE2	1.99	0.44
2:B:17:ILE:HD11	2:B:105:LEU:HD12	1.98	0.44
3:E:63:ASP:H	3:E:66:LYS:HZ3	1.66	0.44
3:D:147:PHE:HB2	3:D:260:ILE:HG23	1.99	0.44
3:D:268:ASP:OD2	3:D:296:ARG:NE	2.51	0.44
3:G:174:TRP:HD1	3:G:191:GLU:HG2	1.83	0.44
1:A:195:PRO:HD2	2:B:26:PRO:HG2	2.00	0.43
3:C:181:ILE:HA	3:C:187:LYS:HZ1	1.83	0.43
3:E:124:THR:O	3:E:128:ASP:HB2	2.19	0.43
3:C:288:VAL:HA	6:M:38:C:H41	1.82	0.43
6:M:36:A:N7	6:M:37:C:N4	2.66	0.43
2:B:147:TRP:CZ2	2:B:153:PRO:HD2	2.54	0.43
3:C:160:VAL:HG13	3:C:213:VAL:HG13	2.01	0.43
3:G:46:GLU:HA	3:G:80:ASP:HA	2.00	0.43
3:H:140:HIS:NE2	3:H:187:LYS:O	2.52	0.43
3:C:301:LYS:HD3	3:C:306:THR:HG21	1.99	0.43
3:D:184:ARG:HE	3:D:279:PRO:HD3	1.84	0.43
3:F:23:ALA:HA	3:F:94:ARG:O	2.18	0.43
3:G:57:LEU:HD11	3:G:67:LEU:HD21	1.99	0.43
3:D:124:THR:O	3:D:128:ASP:CB	2.66	0.43
3:E:16:ARG:NH2	3:F:224:GLU:OE1	2.52	0.42
2:B:240:GLN:HE22	2:B:242:ARG:HH11	1.67	0.42
3:C:137:ARG:HB3	3:C:267:ILE:HD12	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:265:ARG:NH1	3:F:332:ARG:HE	2.17	0.42
3:F:50:ARG:HH21	3:F:74:PRO:HB3	1.84	0.42
3:G:174:TRP:CD1	3:G:191:GLU:HG2	2.54	0.42
3:H:12:LEU:HB2	3:H:336:PHE:HB2	2.01	0.42
3:H:149:TRP:CD1	3:H:259:LYS:HE2	2.55	0.42
3:H:185:ASP:OD1	3:H:185:ASP:N	2.53	0.42
3:D:258:GLN:HB2	6:M:31:C:H5''	2.01	0.42
3:D:184:ARG:NH2	3:D:275:ASP:O	2.52	0.42
3:F:307:LEU:HD22	3:F:318:PRO:HB3	2.02	0.42
3:E:145:ALA:HB2	3:E:158:VAL:HG11	2.01	0.42
3:G:282:VAL:HG12	3:G:329:ASN:HB3	2.02	0.42
3:D:78:THR:HG21	3:D:241:GLN:HE21	1.85	0.42
3:C:200:ALA:HA	3:C:203:LEU:HD12	2.02	0.42
3:D:142:LEU:HD22	3:D:213:VAL:HG21	2.01	0.42
3:H:23:ALA:O	3:H:254:ALA:HA	2.19	0.42
3:D:232:ILE:HD13	3:D:242:LYS:HA	2.01	0.42
3:F:63:ASP:HA	3:F:64:PRO:HD3	1.92	0.42
1:A:233:SER:HB3	2:B:216:LEU:HD11	2.02	0.41
3:C:156:GLU:HB2	3:C:218:ARG:HD3	2.03	0.41
3:F:11:VAL:N	3:F:110:ASN:OD1	2.53	0.41
3:G:265:ARG:NH1	3:G:332:ARG:HE	2.16	0.41
2:B:62:VAL:HG22	2:B:108:SER:H	1.85	0.41
3:D:157:ALA:HB3	3:D:218:ARG:HD3	2.02	0.41
3:G:67:LEU:HD23	3:G:67:LEU:HA	1.94	0.41
2:B:63:CYS:H	2:B:107:VAL:HG23	1.86	0.41
3:E:307:LEU:HD22	3:E:318:PRO:HB3	2.01	0.41
3:G:141:ASN:ND2	3:G:266:THR:O	2.38	0.41
3:C:24:LEU:HD23	3:C:254:ALA:HB2	2.01	0.41
3:D:149:TRP:CE2	3:D:263:ALA:HB2	2.54	0.41
3:C:277:LEU:HD13	3:C:291:GLN:HE22	1.86	0.41
3:D:268:ASP:OD1	3:D:296:ARG:NH2	2.47	0.41
3:H:268:ASP:OD2	3:H:296:ARG:NE	2.44	0.41
3:D:305:TYR:OH	3:E:54:SER:O	2.30	0.41
3:G:136:ARG:O	3:G:140:HIS:ND1	2.54	0.41
3:H:269:THR:HG22	3:H:279:PRO:HB3	2.02	0.41
6:M:30:G:O2'	6:M:31:C:O4'	2.36	0.40
2:B:18:GLN:OE1	2:B:19:ASN:ND2	2.54	0.40
3:D:17:LYS:HG3	3:D:100:LEU:HB2	2.02	0.40
3:G:223:GLN:HG3	3:G:224:GLU:H	1.86	0.40
2:B:10:LEU:HD21	2:B:130:VAL:HG11	2.03	0.40
2:B:308:GLU:HA	2:B:309:PRO:HD3	1.90	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:294:ALA:HB2	3:G:71:ILE:HD13	2.03	0.40
3:D:234:ASP:HB2	3:E:239:LYS:HD3	2.02	0.40
1:A:171:SER:OG	1:A:172:HIS:N	2.55	0.40
2:B:8:ALA:HB3	2:B:113:VAL:HB	2.02	0.40
2:B:199:THR:HG23	2:B:200:LEU:HD12	2.03	0.40
3:E:146:ARG:HG3	3:E:263:ALA:HB1	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	422/436 (97%)	348 (82%)	72 (17%)	2 (0%)	24	57
2	B	301/329 (92%)	259 (86%)	42 (14%)	0	100	100
3	C	287/344 (83%)	260 (91%)	27 (9%)	0	100	100
3	D	331/344 (96%)	292 (88%)	39 (12%)	0	100	100
3	E	332/344 (96%)	304 (92%)	27 (8%)	1 (0%)	36	65
3	F	333/344 (97%)	297 (89%)	36 (11%)	0	100	100
3	G	331/344 (96%)	300 (91%)	31 (9%)	0	100	100
3	H	331/344 (96%)	298 (90%)	33 (10%)	0	100	100
4	K	95/99 (96%)	84 (88%)	11 (12%)	0	100	100
5	L	187/189 (99%)	149 (80%)	38 (20%)	0	100	100
All	All	2950/3117 (95%)	2591 (88%)	356 (12%)	3 (0%)	49	79

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	265	LEU

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Mol	Chain	Res	Type
3	E	250	ARG
1	A	266	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	78/366 (21%)	78 (100%)	0	100	100
2	B	245/271 (90%)	244 (100%)	1 (0%)	84	81
3	C	230/274 (84%)	230 (100%)	0	100	100
3	D	258/274 (94%)	258 (100%)	0	100	100
3	E	259/274 (94%)	258 (100%)	1 (0%)	84	81
3	F	259/274 (94%)	259 (100%)	0	100	100
3	G	259/274 (94%)	259 (100%)	0	100	100
3	H	258/274 (94%)	258 (100%)	0	100	100
4	K	91/86 (106%)	91 (100%)	0	100	100
All	All	1937/2367 (82%)	1935 (100%)	2 (0%)	87	87

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	193	LEU
3	E	33	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	18	GLN
2	B	19	ASN
2	B	121	HIS
2	B	124	GLN
3	C	120	GLN

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Mol	Chain	Res	Type
3	C	223	GLN
3	C	291	GLN
3	C	297	GLN
3	C	324	HIS
3	D	120	GLN
3	D	163	ASN
3	D	223	GLN
3	D	229	GLN
3	D	241	GLN
3	E	120	GLN
3	E	129	GLN
3	E	164	HIS
3	F	120	GLN
3	F	229	GLN
3	F	291	GLN
3	F	324	HIS
3	G	120	GLN
3	G	129	GLN
3	H	129	GLN
3	H	164	HIS
3	H	223	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
6	M	59/60 (98%)	34 (57%)	0

All (34) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
6	M	3	A
6	M	9	U
6	M	10	U
6	M	14	G
6	M	15	G
6	M	16	C
6	M	17	G
6	M	20	C
6	M	21	U
6	M	26	G
6	M	27	U

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Mol	Chain	Res	Type
6	M	28	C
6	M	31	C
6	M	32	G
6	M	33	U
6	M	35	U
6	M	36	A
6	M	37	C
6	M	38	C
6	M	39	U
6	M	40	G
6	M	41	G
6	M	42	U
6	M	43	U
6	M	44	C
6	M	45	A
6	M	46	C
6	M	51	G
6	M	52	U
6	M	53	G
6	M	54	U
6	M	55	A
6	M	59	A
6	M	60	G

There are no RNA pucker outliers to report.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

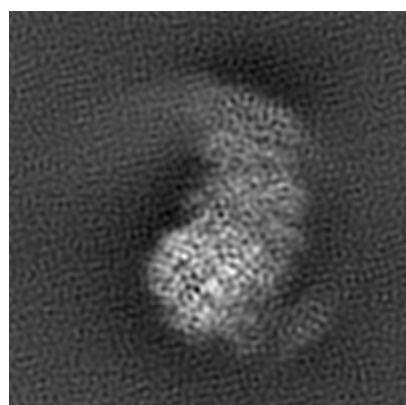
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-7052. These allow visual inspection of the internal detail of the map and identification of artifacts.

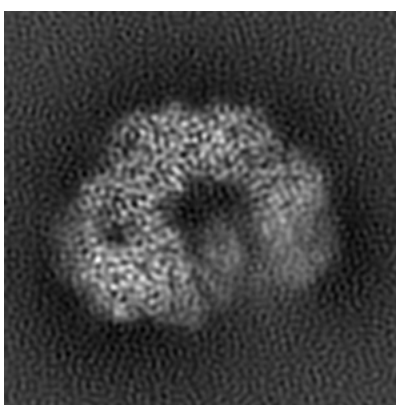
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

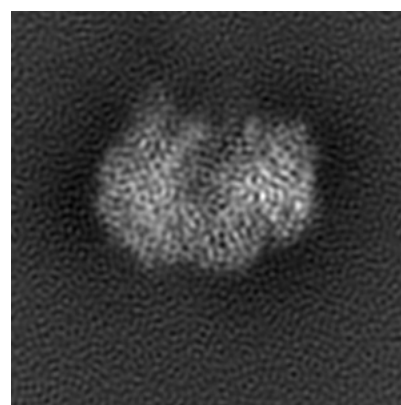
6.1.1 Primary map



X



Y

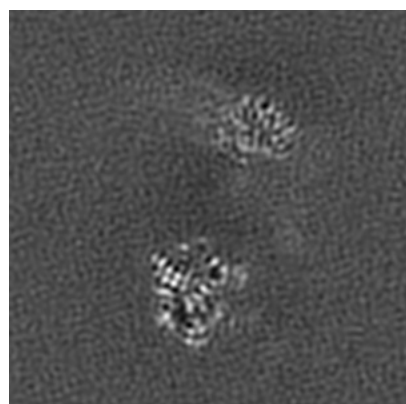


Z

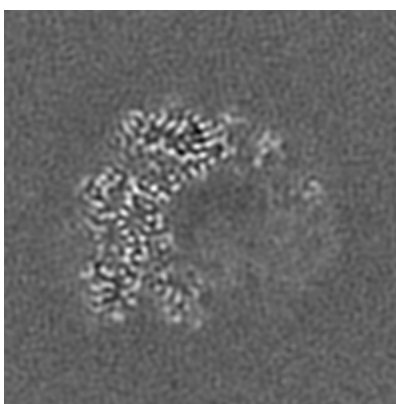
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

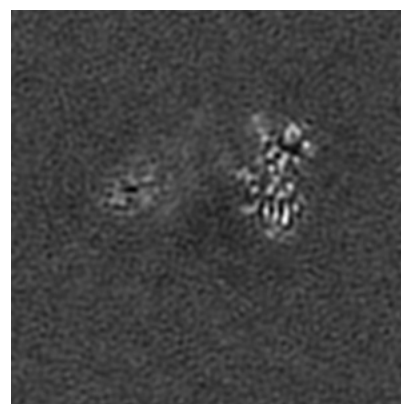
6.2.1 Primary map



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Y Index: 119

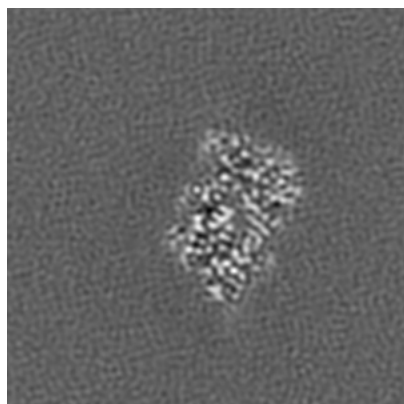


Z Index: 119

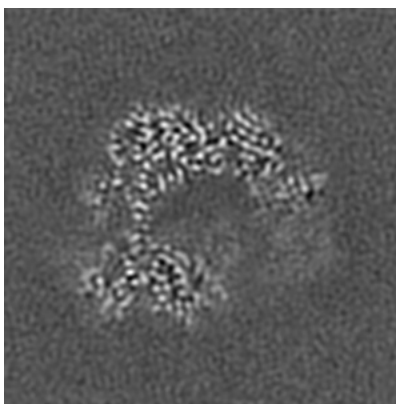
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

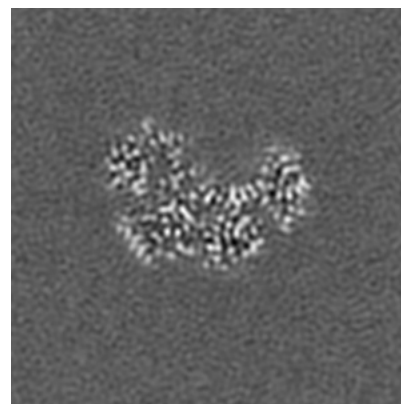
6.3.1 Primary map



X Index: 160



Y Index: 128

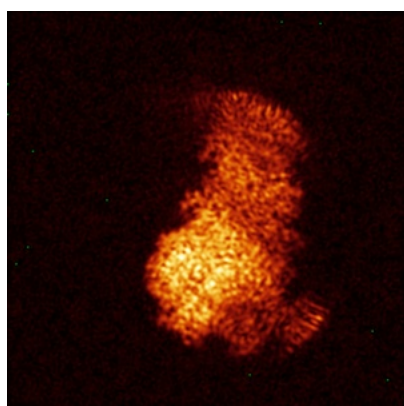


Z Index: 83

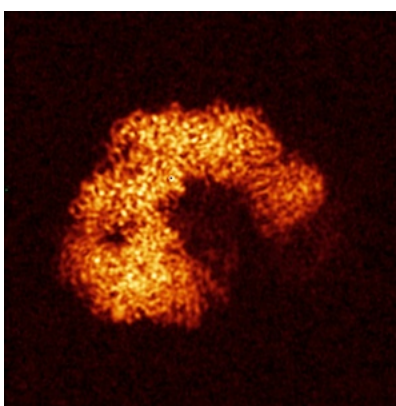
The images above show the largest variance slices of the map in three orthogonal directions.

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

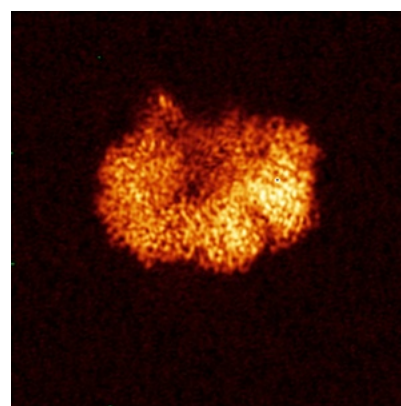
6.4.1 Primary map



X



Y

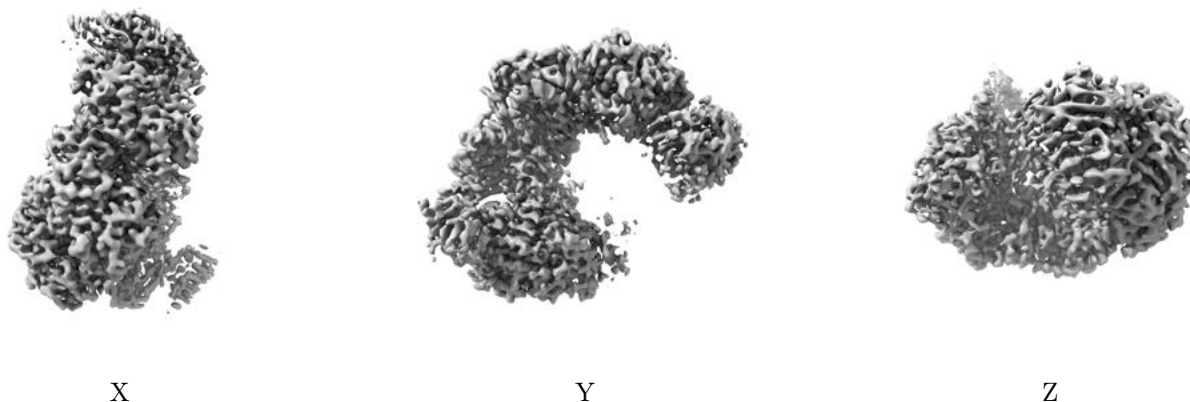


Z

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 2.7. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

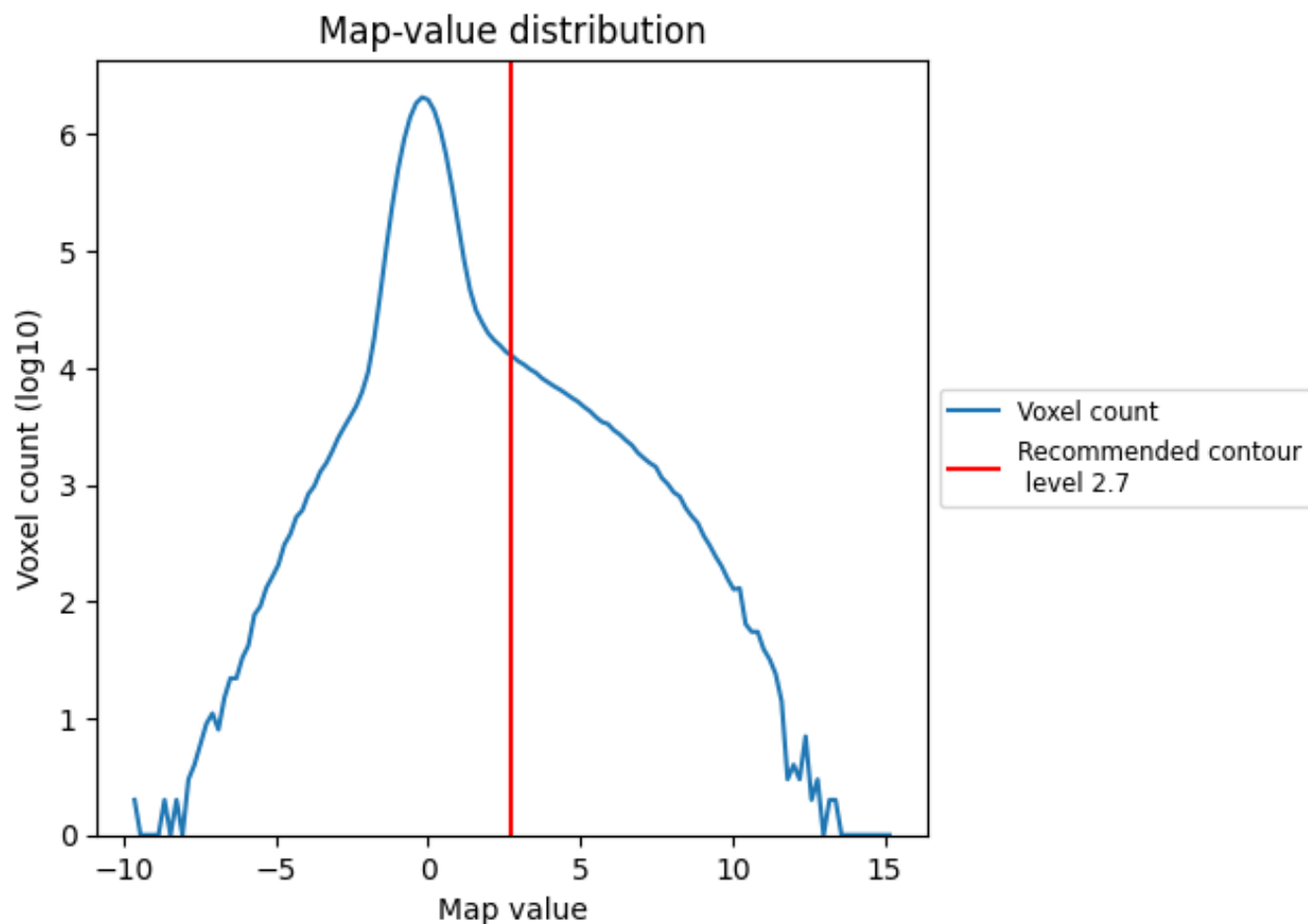
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

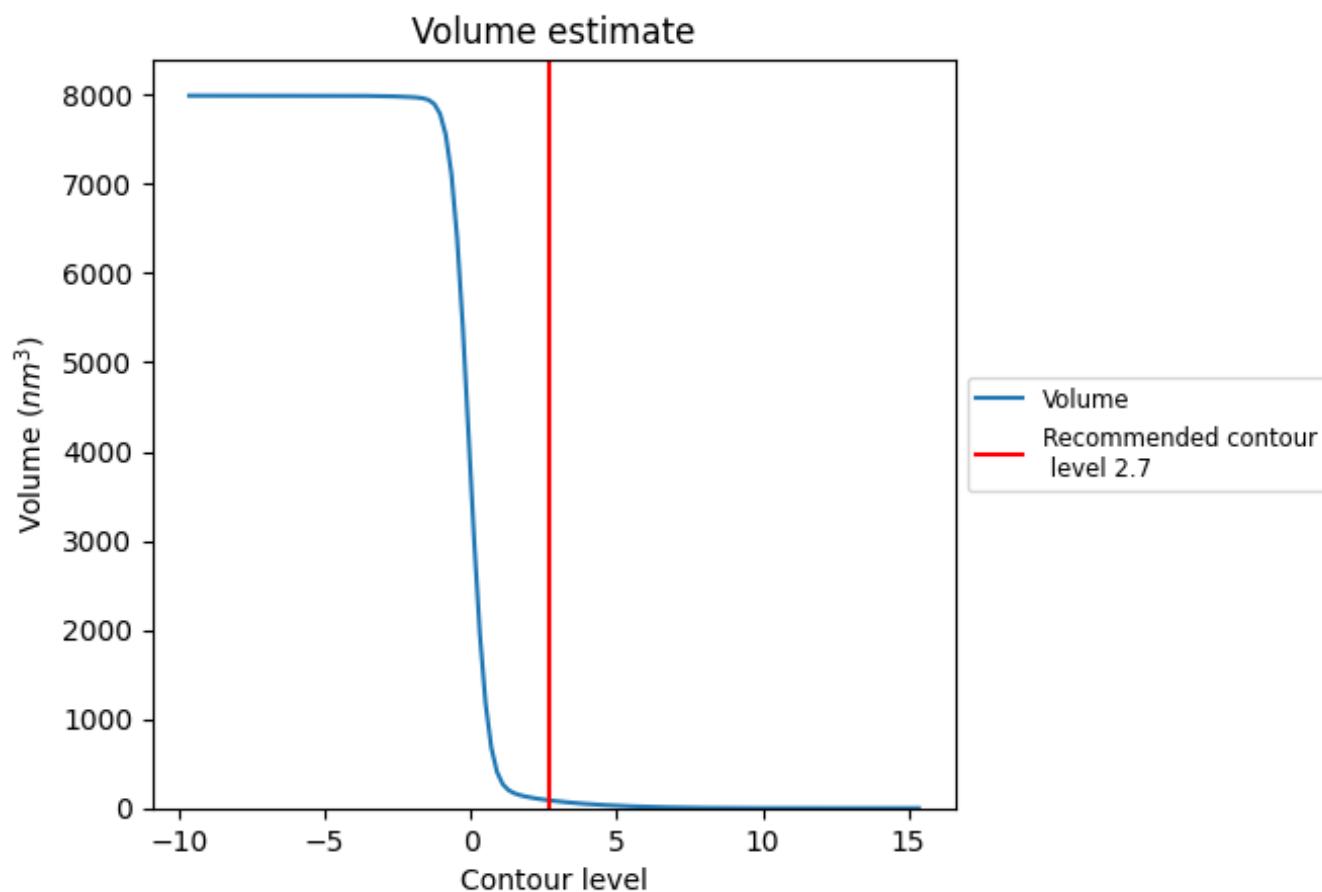
This section contains the results of statistical analysis of the map.

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.

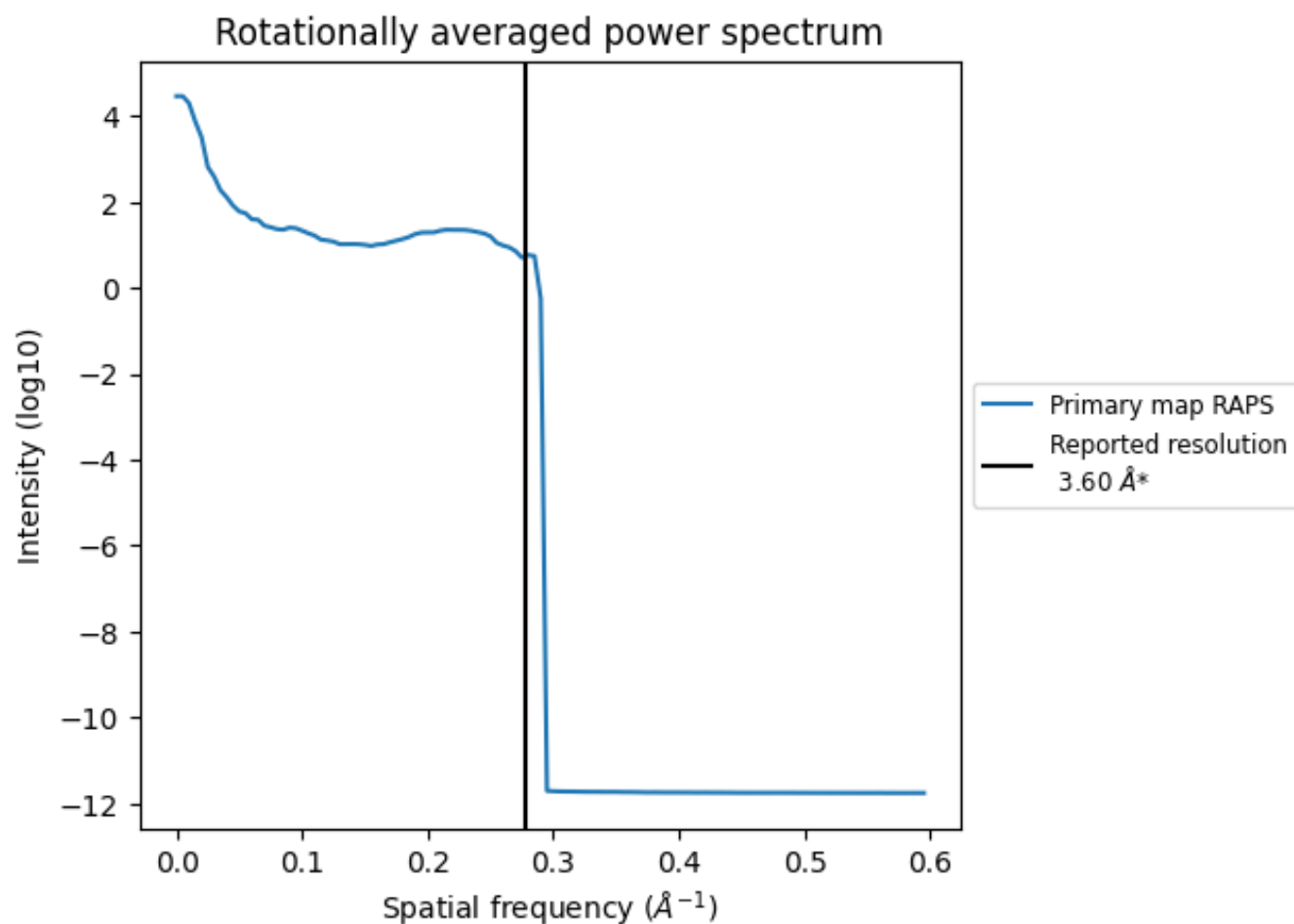
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 87 nm^3 ; this corresponds to an approximate mass of 79 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.278 Å⁻¹

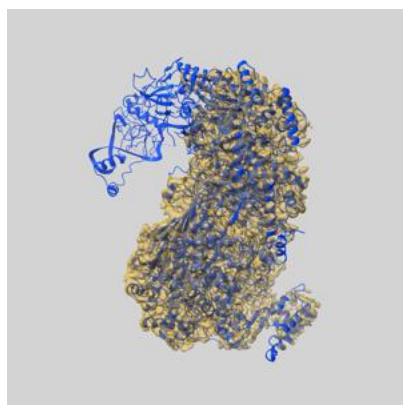
8 Fourier-Shell correlation ⓘ

This section was not generated. No FSC curve or half-maps provided.

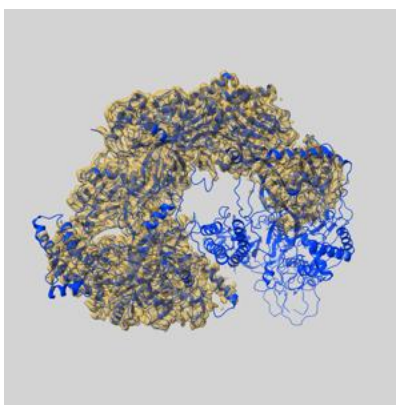
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-7052 and PDB model 6B48. Per-residue inclusion information can be found in section [3](#) on page [7](#).

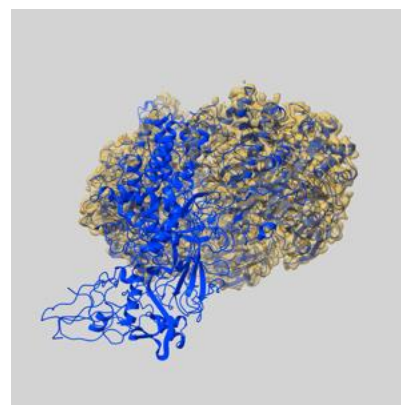
9.1 Map-model overlay [i](#)



X



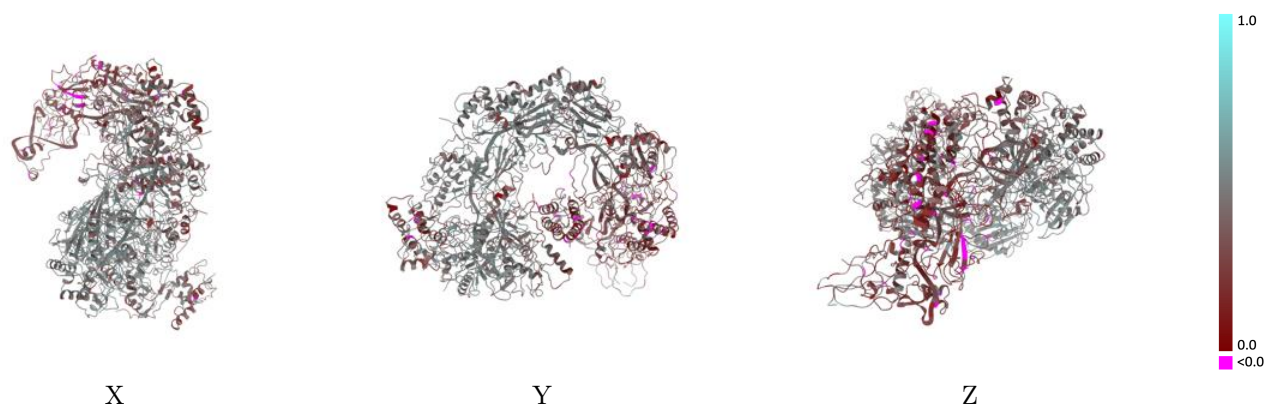
Y



Z

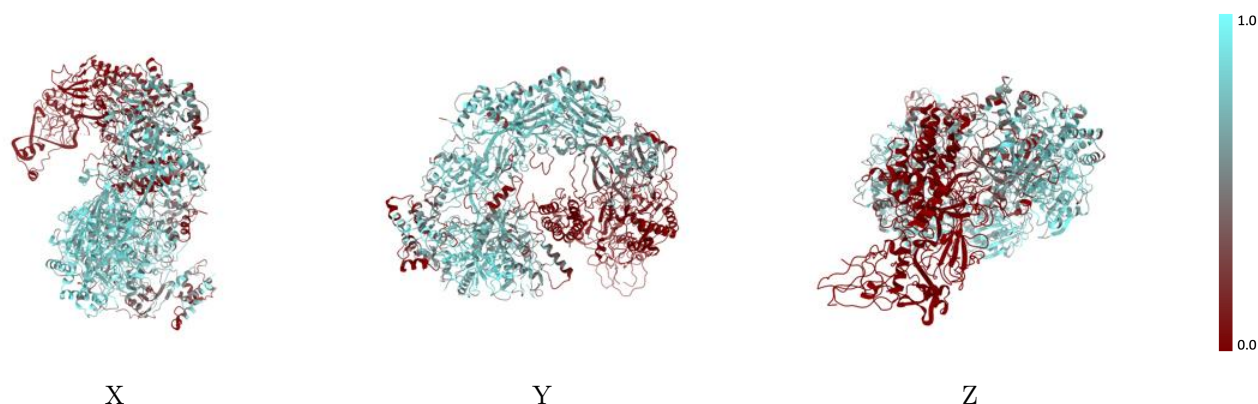
The images above show the 3D surface view of the map at the recommended contour level 2.7 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



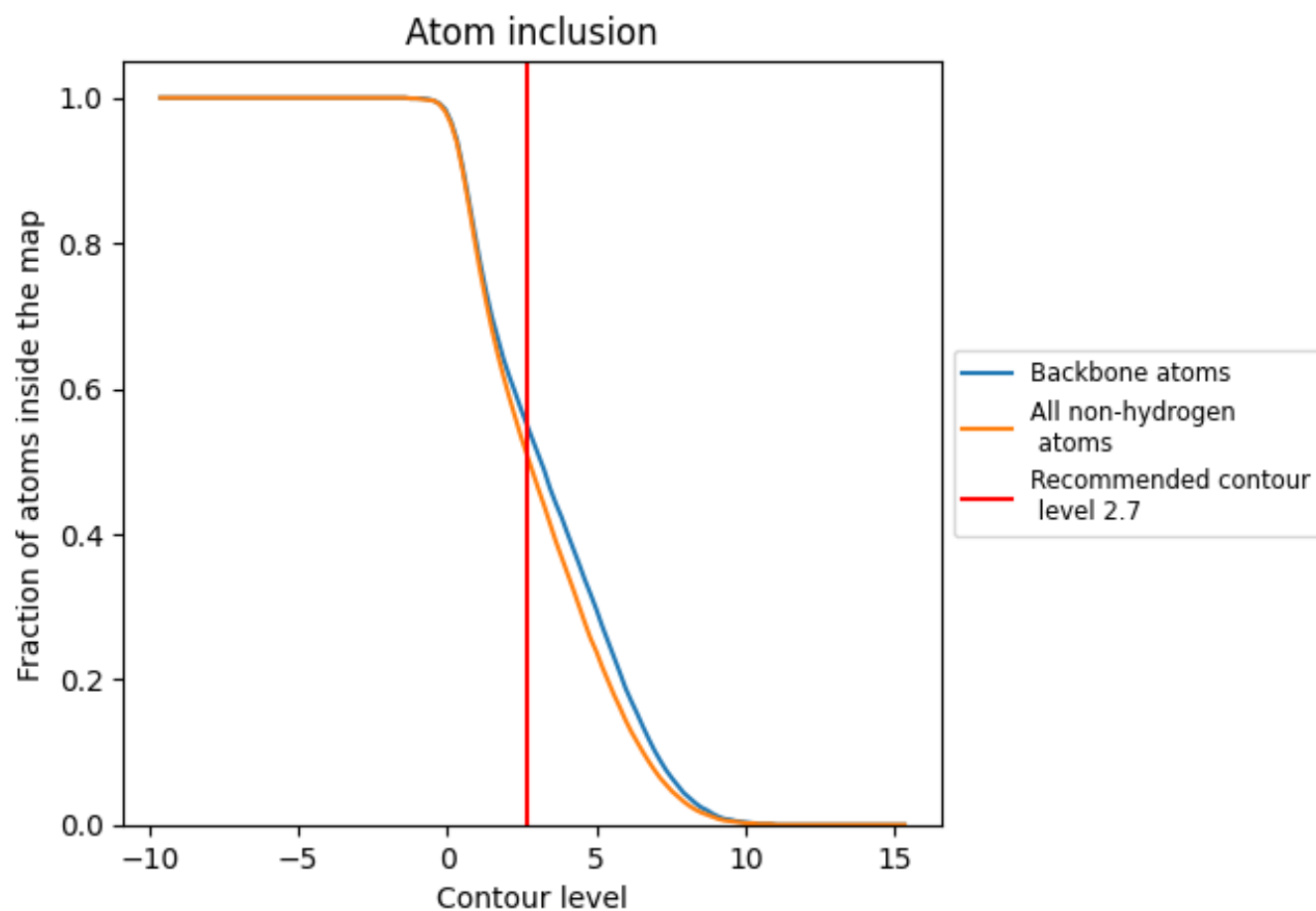
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (2.7).

9.4 Atom inclusion [i](#)



At the recommended contour level, 55% of all backbone atoms, 51% 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 (2.7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5060	 0.4100
A	 0.3930	 0.3860
B	 0.5960	 0.4330
C	 0.0010	 0.2390
D	 0.3430	 0.3490
E	 0.6290	 0.4470
F	 0.7320	 0.4790
G	 0.7510	 0.4830
H	 0.7280	 0.4730
K	 0.4680	 0.4180
L	 0.0000	 0.3180
M	 0.4500	 0.3890

