Full wwPDB EM Validation Report (i)

Feb 26, 2024 – 10:05 PM EST

PDB ID	:	6VQV
EMDB ID	:	EMD-21358
Title	:	Type I-F CRISPR-Csy complex with its inhibitor AcrF9
Authors	:	Zhang, K.; Li, S.; Pintilie, G.; Zhu, Y.; Huang, Z.; Chiu, W.
Deposited on	:	2020-02-06
Resolution	:	2.57 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/EMValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

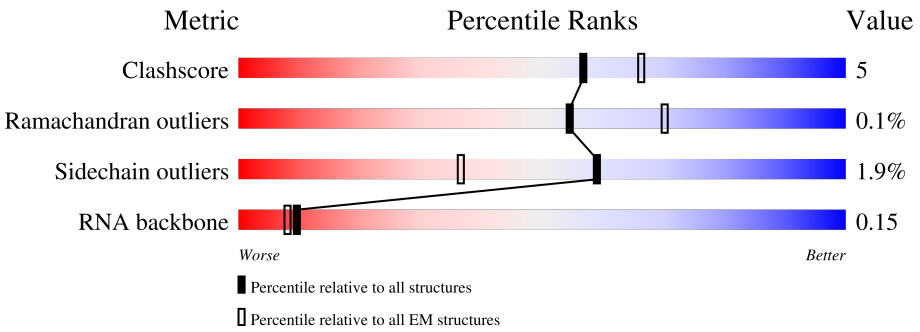
EMDB validation analysis	:	0.0.1.dev70
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.57 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
RNA backbone	4643	859

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion < 40%). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	68	9% 84% 12% • •
1	B	68	26% 81% 15% • •
2	C	434	62% 82% 15% •
3	D	327	13% 79% 15% 6%
4	E	360	25% 64% 15% 21%
4	F	360	11% 79% 12% 9%
4	G	360	6% 79% 10% • 9%

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Mol	Chain	Length	Quality of chain
4	H	360	6% 81% 11% • 7%
4	I	360	5% 76% 14% 9%
4	J	360	9% 82% 11% • 6%
5	K	187	94% 84% 14% ••
6	L	60	35% 37% 50% 13%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 24676 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 AcrF9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	68	545	339	98	104	4	0	0
1	B	68	545	339	98	104	4	0	0

- Molecule 2 is a protein called CRISPR-associated protein Csy1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	C	422	3378	2116	636	622	4	0	0

- Molecule 3 is a protein called CRISPR-associated protein Csy2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	D	306	2410	1527	445	433	5	0	0

- Molecule 4 is a protein called CRISPR-associated protein Csy3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	E	285	2204	1390	396	416	2	0	0
4	F	328	2541	1595	462	482	2	0	0
4	G	327	2532	1589	460	481	2	0	0
4	H	335	2592	1626	471	493	2	0	0
4	I	327	2532	1589	460	481	2	0	0
4	J	337	2611	1636	474	499	2	0	0

There are 114 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	1	MET	-	expression tag	UNP A0A444M080
E	2	LYS	-	expression tag	UNP A0A444M080
E	3	SER	-	expression tag	UNP A0A444M080
E	4	SER	-	expression tag	UNP A0A444M080
E	5	HIS	-	expression tag	UNP A0A444M080
E	6	HIS	-	expression tag	UNP A0A444M080
E	7	HIS	-	expression tag	UNP A0A444M080
E	8	HIS	-	expression tag	UNP A0A444M080
E	9	HIS	-	expression tag	UNP A0A444M080
E	10	HIS	-	expression tag	UNP A0A444M080
E	11	GLU	-	expression tag	UNP A0A444M080
E	12	ASN	-	expression tag	UNP A0A444M080
E	13	LEU	-	expression tag	UNP A0A444M080
E	14	TYR	-	expression tag	UNP A0A444M080
E	15	PHE	-	expression tag	UNP A0A444M080
E	16	GLN	-	expression tag	UNP A0A444M080
E	17	SER	-	expression tag	UNP A0A444M080
E	18	ASN	-	expression tag	UNP A0A444M080
E	19	ALA	-	expression tag	UNP A0A444M080
F	1	MET	-	expression tag	UNP A0A444M080
F	2	LYS	-	expression tag	UNP A0A444M080
F	3	SER	-	expression tag	UNP A0A444M080
F	4	SER	-	expression tag	UNP A0A444M080
F	5	HIS	-	expression tag	UNP A0A444M080
F	6	HIS	-	expression tag	UNP A0A444M080
F	7	HIS	-	expression tag	UNP A0A444M080
F	8	HIS	-	expression tag	UNP A0A444M080
F	9	HIS	-	expression tag	UNP A0A444M080
F	10	HIS	-	expression tag	UNP A0A444M080
F	11	GLU	-	expression tag	UNP A0A444M080
F	12	ASN	-	expression tag	UNP A0A444M080
F	13	LEU	-	expression tag	UNP A0A444M080
F	14	TYR	-	expression tag	UNP A0A444M080
F	15	PHE	-	expression tag	UNP A0A444M080
F	16	GLN	-	expression tag	UNP A0A444M080
F	17	SER	-	expression tag	UNP A0A444M080
F	18	ASN	-	expression tag	UNP A0A444M080
F	19	ALA	-	expression tag	UNP A0A444M080
G	1	MET	-	expression tag	UNP A0A444M080
G	2	LYS	-	expression tag	UNP A0A444M080
G	3	SER	-	expression tag	UNP A0A444M080
G	4	SER	-	expression tag	UNP A0A444M080

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Chain	Residue	Modelled	Actual	Comment	Reference
G	5	HIS	-	expression tag	UNP A0A444M080
G	6	HIS	-	expression tag	UNP A0A444M080
G	7	HIS	-	expression tag	UNP A0A444M080
G	8	HIS	-	expression tag	UNP A0A444M080
G	9	HIS	-	expression tag	UNP A0A444M080
G	10	HIS	-	expression tag	UNP A0A444M080
G	11	GLU	-	expression tag	UNP A0A444M080
G	12	ASN	-	expression tag	UNP A0A444M080
G	13	LEU	-	expression tag	UNP A0A444M080
G	14	TYR	-	expression tag	UNP A0A444M080
G	15	PHE	-	expression tag	UNP A0A444M080
G	16	GLN	-	expression tag	UNP A0A444M080
G	17	SER	-	expression tag	UNP A0A444M080
G	18	ASN	-	expression tag	UNP A0A444M080
G	19	ALA	-	expression tag	UNP A0A444M080
H	1	MET	-	expression tag	UNP A0A444M080
H	2	LYS	-	expression tag	UNP A0A444M080
H	3	SER	-	expression tag	UNP A0A444M080
H	4	SER	-	expression tag	UNP A0A444M080
H	5	HIS	-	expression tag	UNP A0A444M080
H	6	HIS	-	expression tag	UNP A0A444M080
H	7	HIS	-	expression tag	UNP A0A444M080
H	8	HIS	-	expression tag	UNP A0A444M080
H	9	HIS	-	expression tag	UNP A0A444M080
H	10	HIS	-	expression tag	UNP A0A444M080
H	11	GLU	-	expression tag	UNP A0A444M080
H	12	ASN	-	expression tag	UNP A0A444M080
H	13	LEU	-	expression tag	UNP A0A444M080
H	14	TYR	-	expression tag	UNP A0A444M080
H	15	PHE	-	expression tag	UNP A0A444M080
H	16	GLN	-	expression tag	UNP A0A444M080
H	17	SER	-	expression tag	UNP A0A444M080
H	18	ASN	-	expression tag	UNP A0A444M080
H	19	ALA	-	expression tag	UNP A0A444M080
I	1	MET	-	expression tag	UNP A0A444M080
I	2	LYS	-	expression tag	UNP A0A444M080
I	3	SER	-	expression tag	UNP A0A444M080
I	4	SER	-	expression tag	UNP A0A444M080
I	5	HIS	-	expression tag	UNP A0A444M080
I	6	HIS	-	expression tag	UNP A0A444M080
I	7	HIS	-	expression tag	UNP A0A444M080
I	8	HIS	-	expression tag	UNP A0A444M080

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Chain	Residue	Modelled	Actual	Comment	Reference
I	9	HIS	-	expression tag	UNP A0A444M080
I	10	HIS	-	expression tag	UNP A0A444M080
I	11	GLU	-	expression tag	UNP A0A444M080
I	12	ASN	-	expression tag	UNP A0A444M080
I	13	LEU	-	expression tag	UNP A0A444M080
I	14	TYR	-	expression tag	UNP A0A444M080
I	15	PHE	-	expression tag	UNP A0A444M080
I	16	GLN	-	expression tag	UNP A0A444M080
I	17	SER	-	expression tag	UNP A0A444M080
I	18	ASN	-	expression tag	UNP A0A444M080
I	19	ALA	-	expression tag	UNP A0A444M080
J	1	MET	-	expression tag	UNP A0A444M080
J	2	LYS	-	expression tag	UNP A0A444M080
J	3	SER	-	expression tag	UNP A0A444M080
J	4	SER	-	expression tag	UNP A0A444M080
J	5	HIS	-	expression tag	UNP A0A444M080
J	6	HIS	-	expression tag	UNP A0A444M080
J	7	HIS	-	expression tag	UNP A0A444M080
J	8	HIS	-	expression tag	UNP A0A444M080
J	9	HIS	-	expression tag	UNP A0A444M080
J	10	HIS	-	expression tag	UNP A0A444M080
J	11	GLU	-	expression tag	UNP A0A444M080
J	12	ASN	-	expression tag	UNP A0A444M080
J	13	LEU	-	expression tag	UNP A0A444M080
J	14	TYR	-	expression tag	UNP A0A444M080
J	15	PHE	-	expression tag	UNP A0A444M080
J	16	GLN	-	expression tag	UNP A0A444M080
J	17	SER	-	expression tag	UNP A0A444M080
J	18	ASN	-	expression tag	UNP A0A444M080
J	19	ALA	-	expression tag	UNP A0A444M080

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	K	187	1511	953	291	263	4	0	0

- Molecule 6 is a RNA chain called CrRNA (60-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
6	L	60	1275	569	223	423	60	0	0

There is a discrepancy between the modelled and reference sequences:

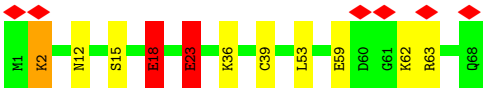
Chain	Residue	Modelled	Actual	Comment	Reference
L	53	A	G	conflict	GB 313291946

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.

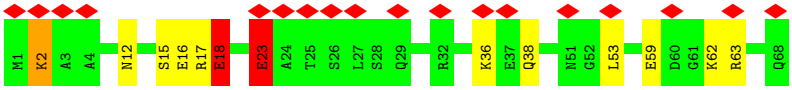
- Molecule 1: AcrF9

Chain A:



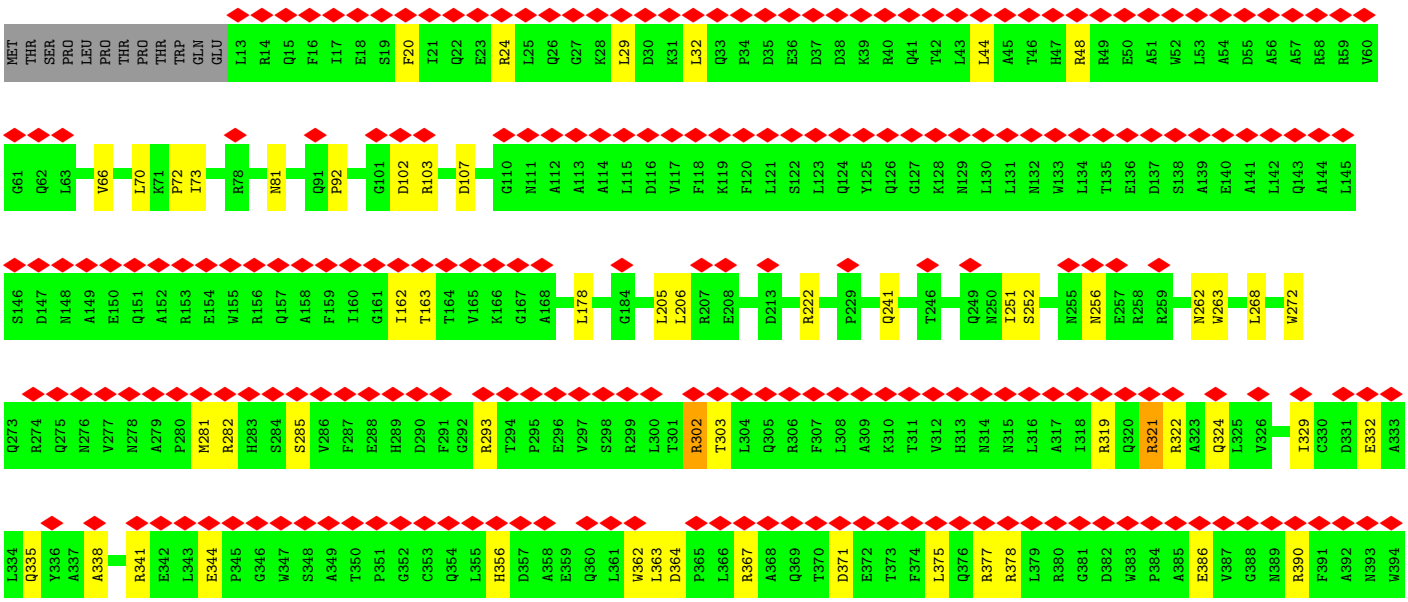
- Molecule 1: AcrF9

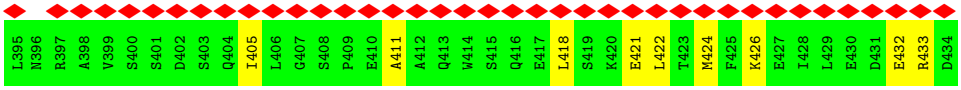
Chain B:



- Molecule 2: CRISPR-associated protein Csy1

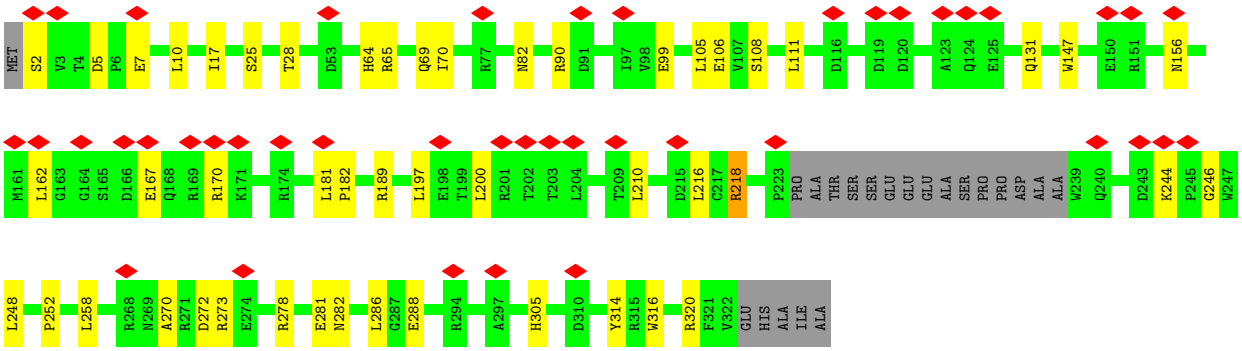
Chain C:





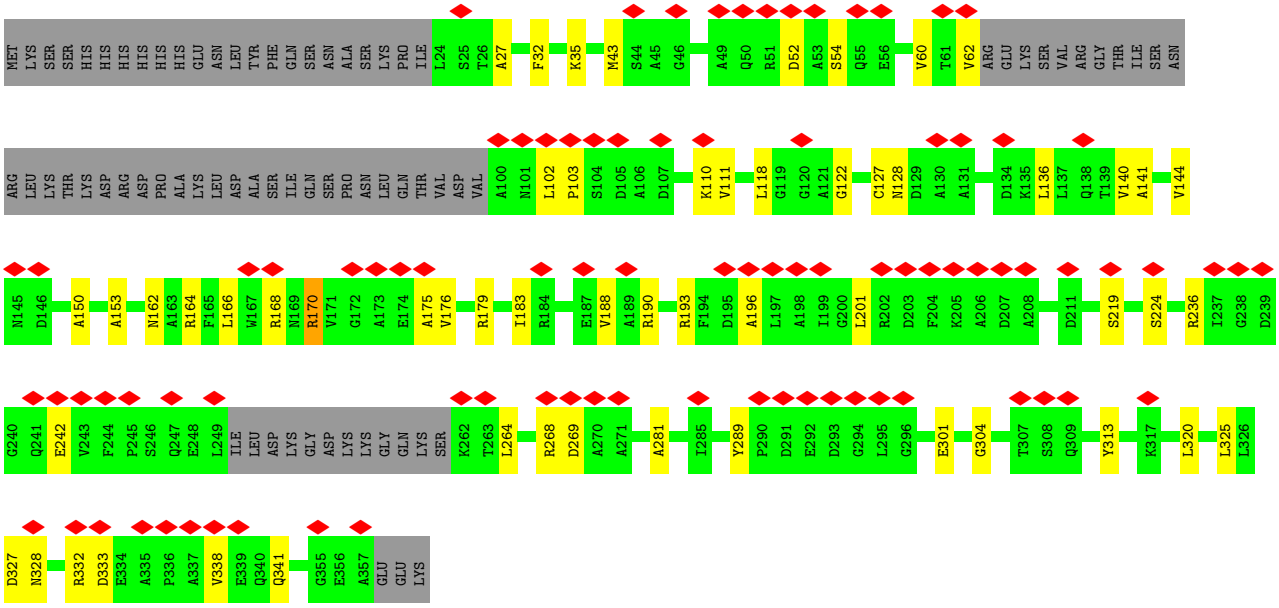
• Molecule 3: CRISPR-associated protein Csy2

Chain D:



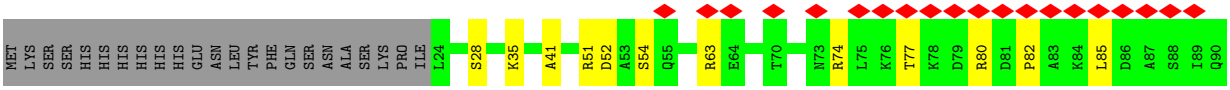
• Molecule 4: CRISPR-associated protein Csy3

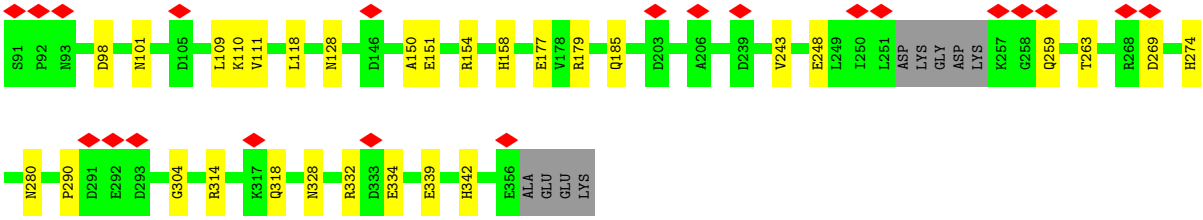
Chain E:



• Molecule 4: CRISPR-associated protein Csy3

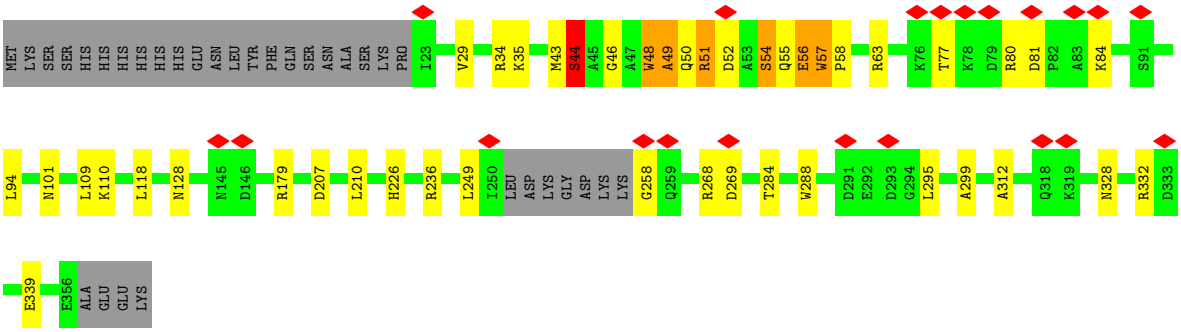
Chain F:





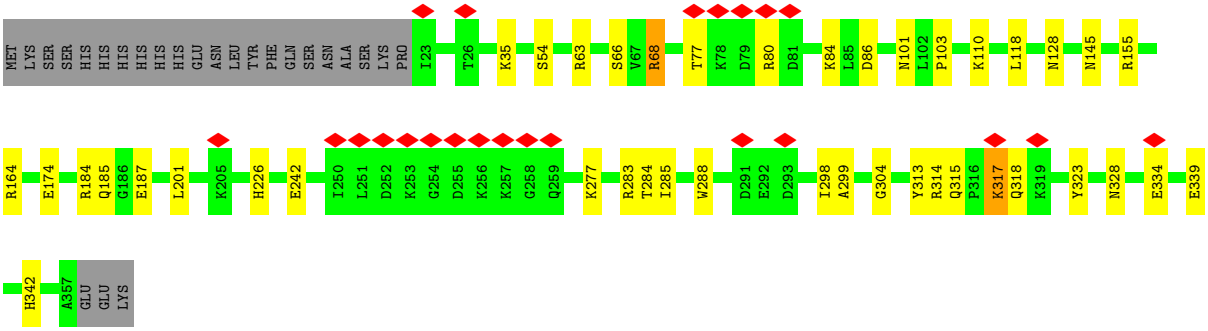
• Molecule 4: CRISPR-associated protein Csy3

Chain G:



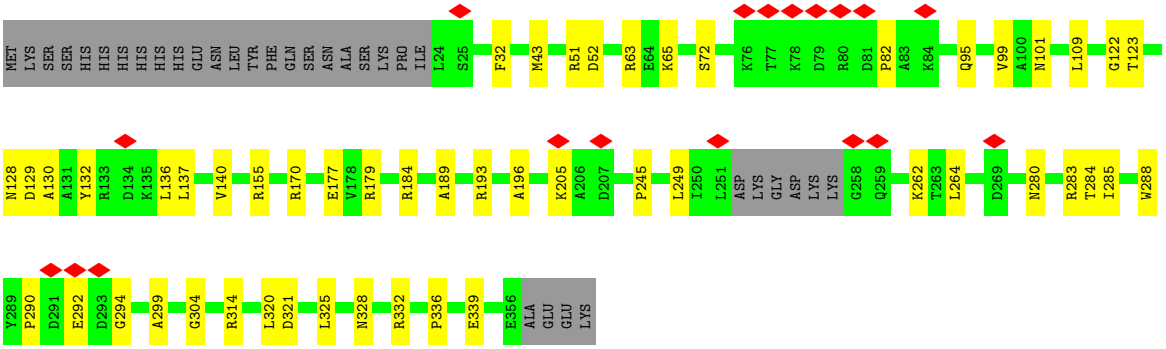
• Molecule 4: CRISPR-associated protein Csy3

Chain H:



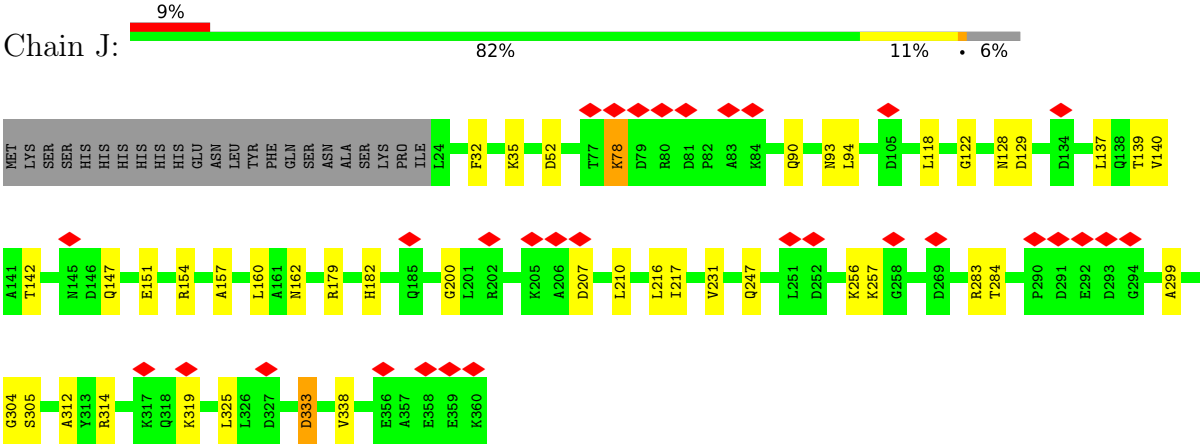
• Molecule 4: CRISPR-associated protein Csy3

Chain I:



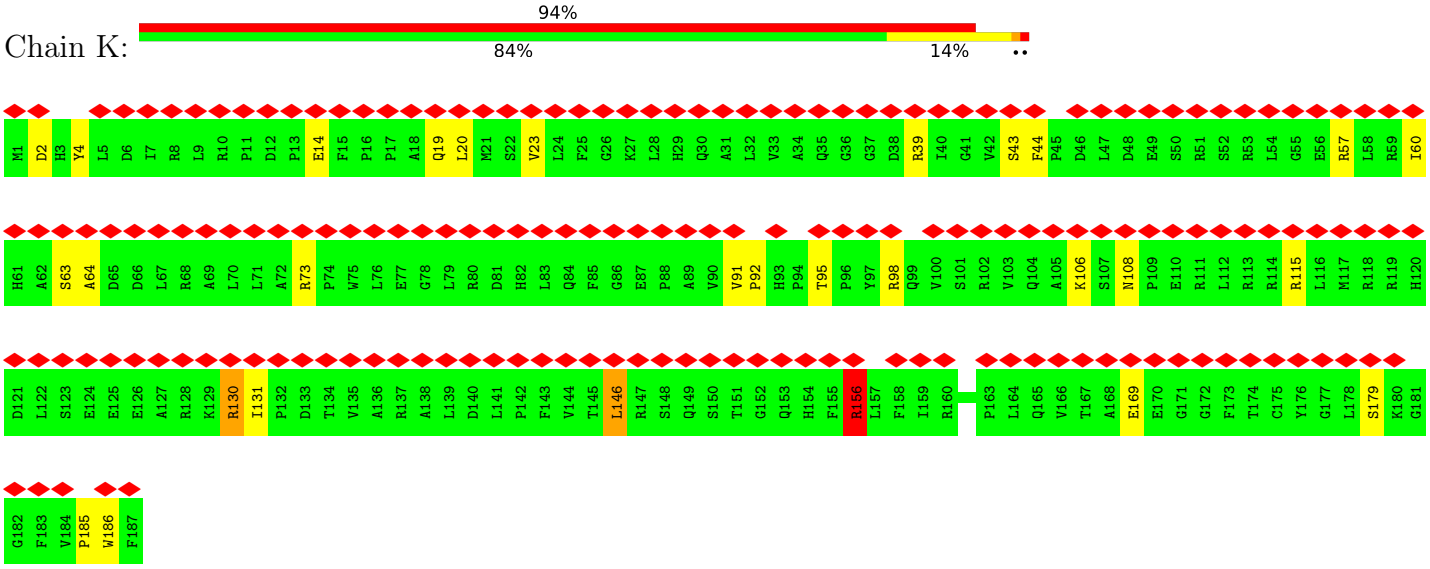
- Molecule 4: CRISPR-associated protein Csy3

Chain J:



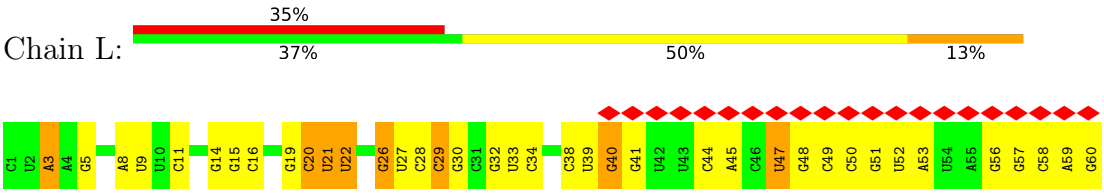
- Molecule 5: CRISPR-associated endonuclease Cas6/Csy4

Chain K:



- Molecule 6: CrRNA (60-MER)

Chain L:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	332404	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 ($e^-/Å^2$)	7	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	3.471	Depositor
Minimum map value	-1.533	Depositor
Average map value	0.006	Depositor
Map value standard deviation	0.089	Depositor
Recommended contour level	0.43	Depositor
Map size (Å)	249.59999, 249.59999, 249.59999	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.65, 0.65, 0.65	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	0.49	0/552	0.76	2/740 (0.3%)
1	B	0.49	0/552	0.76	2/740 (0.3%)
2	C	0.35	0/3459	0.60	1/4694 (0.0%)
3	D	0.41	0/2469	0.64	0/3361
4	E	0.37	0/2247	0.66	2/3052 (0.1%)
4	F	0.43	0/2587	0.59	1/3509 (0.0%)
4	G	0.73	5/2578 (0.2%)	0.77	9/3498 (0.3%)
4	H	0.49	0/2639	0.62	0/3579
4	I	0.50	0/2578	0.59	0/3498
4	J	0.47	0/2658	0.64	2/3603 (0.1%)
5	K	0.38	1/1550 (0.1%)	0.64	1/2096 (0.0%)
6	L	0.82	0/1423	1.09	7/2215 (0.3%)
All	All	0.50	6/25292 (0.0%)	0.68	27/34585 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
3	D	0	1
4	E	0	1
5	K	0	1
All	All	0	3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	57	TRP	CG-CD1	-8.69	1.24	1.36
4	G	57	TRP	CB-CG	-7.62	1.36	1.50
4	G	46	GLY	C-O	-7.35	1.11	1.23
5	K	186	TRP	CD2-CE2	-6.66	1.33	1.41
4	G	58	PRO	CA-C	-5.72	1.41	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	44	SER	C-O	-5.69	1.12	1.23

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	L	29	C	C6-N1-C2	-12.16	115.44	120.30
6	L	29	C	C5-C6-N1	9.90	125.95	121.00
4	G	269	ASP	CB-CG-OD1	8.63	126.07	118.30
6	L	29	C	N1-C2-O2	8.15	123.79	118.90
6	L	29	C	N3-C2-O2	-7.64	116.55	121.90
4	G	57	TRP	CB-CA-C	7.27	124.94	110.40
1	A	18	GLU	CA-CB-CG	7.23	129.31	113.40
1	B	18	GLU	CA-CB-CG	7.23	129.30	113.40
4	G	44	SER	CB-CA-C	-6.79	97.19	110.10
4	G	48	TRP	C-N-CA	6.79	138.67	121.70
4	G	57	TRP	CA-CB-CG	6.63	126.30	113.70
4	G	295	LEU	CA-CB-CG	6.57	130.42	115.30
4	G	48	TRP	CA-CB-CG	6.54	126.13	113.70
4	G	54	SER	N-CA-CB	6.50	120.26	110.50
6	L	29	C	C2-N1-C1'	6.44	125.88	118.80
4	J	333	ASP	CB-CG-OD1	6.42	124.08	118.30
6	L	21	U	C2-N1-C1'	6.32	125.28	117.70
4	E	269	ASP	CB-CG-OD1	5.95	123.65	118.30
5	K	146	LEU	CA-CB-CG	5.86	128.78	115.30
2	C	205	LEU	CA-CB-CG	5.62	128.23	115.30
6	L	21	U	N3-C2-O2	-5.62	118.27	122.20
4	F	269	ASP	CB-CG-OD1	5.55	123.29	118.30
4	G	49	ALA	N-CA-C	-5.51	96.11	111.00
1	B	23	GLU	CA-CB-CG	5.46	125.41	113.40
1	A	23	GLU	CA-CB-CG	5.43	125.34	113.40
4	E	333	ASP	CB-CG-OD1	5.26	123.03	118.30
4	J	256	LYS	C-N-CA	5.05	134.32	121.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	D	147	TRP	Peptide
4	E	268	ARG	Peptide
5	K	156	ARG	Peptide

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	545	0	547	6	0
1	B	545	0	547	8	0
2	C	3378	0	3311	33	0
3	D	2410	0	2406	29	0
4	E	2204	0	2163	31	0
4	F	2541	0	2527	23	0
4	G	2532	0	2514	40	0
4	H	2592	0	2581	29	0
4	I	2532	0	2514	29	0
4	J	2611	0	2595	26	0
5	K	1511	0	1510	15	0
6	L	1275	0	646	12	0
All	All	24676	0	23861	248	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 5.

All (248) 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)
4:G:51:ARG:NH1	4:G:52:ASP:HB2	1.66	1.09
4:G:51:ARG:HH11	4:G:52:ASP:HB2	1.09	1.08
4:G:51:ARG:HB3	4:G:179:ARG:HH22	1.42	0.85
4:G:51:ARG:HH11	4:G:52:ASP:CB	1.94	0.79
4:G:48:TRP:CD1	4:G:236:ARG:NH2	2.52	0.78
4:G:48:TRP:CG	4:G:236:ARG:NH2	2.53	0.76
4:G:49:ALA:HA	4:G:51:ARG:NH2	2.01	0.74
4:G:51:ARG:CB	4:G:179:ARG:HH22	2.03	0.72
4:G:54:SER:HB2	4:G:110:LYS:HD2	1.71	0.71
2:C:20:PHE:HB3	2:C:24:ARG:HH12	1.55	0.70
4:G:49:ALA:HA	4:G:51:ARG:HH22	1.59	0.68
4:G:51:ARG:HB3	4:G:179:ARG:NH2	2.09	0.68
4:E:111:VAL:HG21	4:E:166:LEU:HD21	1.75	0.68
4:G:51:ARG:NH1	4:G:52:ASP:CB	2.50	0.67
4:H:288:TRP:O	4:H:314:ARG:NH2	2.30	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:48:TRP:CZ3	4:G:236:ARG:HG3	2.33	0.63
1:B:38:GLN:O	4:J:78:LYS:NZ	2.30	0.63
3:D:7:GLU:HA	3:D:162:LEU:HD13	1.81	0.63
4:G:51:ARG:HA	4:G:179:ARG:HH12	1.63	0.63
4:E:175:ALA:HB3	4:E:236:ARG:HD3	1.79	0.62
3:D:270:ALA:O	4:J:128:ASN:ND2	2.33	0.62
2:C:371:ASP:H	2:C:378:ARG:HD2	1.64	0.61
4:H:185:GLN:HB2	4:H:187:GLU:HG3	1.83	0.61
3:D:273:ARG:NH1	4:J:129:ASP:OD1	2.34	0.61
1:A:18:GLU:N	1:A:18:GLU:OE1	2.34	0.60
4:E:62:VAL:HG21	4:E:264:LEU:HD11	1.84	0.60
4:E:60:VAL:HG13	4:E:102:LEU:HD21	1.83	0.60
4:I:65:LYS:HE3	4:I:99:VAL:HG11	1.83	0.60
1:B:59:GLU:HG3	1:B:62:LYS:HD2	1.83	0.60
4:I:184:ARG:HG2	4:I:189:ALA:HB2	1.83	0.60
1:A:59:GLU:HG3	1:A:62:LYS:HD2	1.83	0.59
2:C:364:ASP:O	2:C:433:ARG:NH2	2.35	0.59
1:B:17:ARG:HD2	4:J:93:ASN:HB3	1.83	0.59
4:H:63:ARG:NH2	4:H:101:ASN:OD1	2.36	0.58
4:I:51:ARG:NH2	4:I:177:GLU:OE1	2.36	0.58
4:I:290:PRO:HB3	4:I:320:LEU:HD12	1.85	0.58
4:G:54:SER:HB2	4:G:110:LYS:CD	2.33	0.57
4:E:328:ASN:HD22	4:E:332:ARG:HD2	1.69	0.57
4:E:164:ARG:HG3	4:E:281:ALA:HB1	1.86	0.57
2:C:70:LEU:HD12	2:C:73:ILE:HD12	1.87	0.57
4:F:98:ASP:OD2	4:F:259:GLN:NE2	2.38	0.57
4:J:319:LYS:HZ2	4:J:325:LEU:HD23	1.70	0.57
3:D:189:ARG:NH1	3:D:288:GLU:OE1	2.38	0.56
5:K:19:GLN:HG3	5:K:20:LEU:HG	1.85	0.56
2:C:321:ARG:HA	2:C:324:GLN:HG2	1.88	0.56
4:F:54:SER:O	4:F:110:LYS:NZ	2.39	0.55
1:B:18:GLU:OE1	1:B:18:GLU:N	2.34	0.55
2:C:281:MET:HG3	2:C:282:ARG:HG3	1.88	0.55
4:H:155:ARG:HG2	4:H:285:ILE:HG23	1.88	0.55
4:E:179:ARG:HG2	4:E:193:ARG:HG2	1.88	0.55
4:I:304:GLY:H	4:I:314:ARG:HB2	1.72	0.55
4:J:162:ASN:ND2	4:J:200:GLY:O	2.38	0.55
4:E:122:GLY:HA3	4:E:140:VAL:HG21	1.89	0.55
4:H:35:LYS:HG3	4:H:118:LEU:HB2	1.89	0.54
4:F:304:GLY:H	4:F:314:ARG:HB2	1.71	0.54
4:G:51:ARG:CA	4:G:179:ARG:HH22	2.20	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:184:ARG:NH1	4:H:226:HIS:O	2.41	0.54
4:F:328:ASN:HB3	4:F:334:GLU:HB3	1.89	0.54
4:I:284:THR:HA	4:I:299:ALA:HA	1.88	0.54
4:E:183:ILE:HG22	4:E:188:VAL:HA	1.89	0.54
4:I:288:TRP:O	4:I:314:ARG:NH2	2.40	0.54
4:H:128:ASN:ND2	4:J:90:GLN:OE1	2.40	0.54
5:K:108:ASN:H	5:K:156:ARG:HH22	1.55	0.54
4:H:283:ARG:NH2	6:L:20:C:OP2	2.41	0.54
4:J:283:ARG:NH2	6:L:8:A:OP2	2.38	0.54
4:G:56:GLU:O	4:G:56:GLU:HG2	2.07	0.54
2:C:432:GLU:HG2	2:C:433:ARG:HG2	1.90	0.53
5:K:44:PHE:O	5:K:98:ARG:NH1	2.41	0.53
4:J:147:GLN:HE22	4:J:338:VAL:HG13	1.73	0.53
5:K:23:VAL:HG22	5:K:146:LEU:HA	1.90	0.53
5:K:115:ARG:HH12	6:L:47:U:H4'	1.74	0.53
4:E:127:CYS:SG	4:E:128:ASN:N	2.82	0.53
4:F:51:ARG:NE	4:F:177:GLU:OE1	2.34	0.52
4:H:315:GLN:H	4:H:318:GLN:HE21	1.56	0.52
3:D:244:LYS:HG2	3:D:246:GLY:H	1.75	0.52
4:I:328:ASN:OD1	4:I:332:ARG:NH1	2.42	0.52
5:K:43:SER:H	5:K:60:ILE:HG23	1.75	0.52
2:C:162:ILE:HG23	2:C:163:THR:HG23	1.91	0.52
4:G:35:LYS:HG3	4:G:118:LEU:HB2	1.92	0.52
2:C:107:ASP:OD1	2:C:252:SER:OG	2.27	0.52
3:D:5:ASP:HA	3:D:320:ARG:HH12	1.75	0.52
3:D:17:ILE:HB	3:D:105:LEU:HB3	1.93	0.51
2:C:178:LEU:HD11	6:L:3:A:H5"	1.92	0.51
2:C:344:GLU:O	2:C:367:ARG:NH2	2.43	0.51
4:I:292:GLU:HG2	4:I:294:GLY:H	1.76	0.51
4:F:150:ALA:O	4:F:154:ARG:HB2	2.11	0.51
2:C:422:LEU:O	2:C:426:LYS:NZ	2.37	0.51
3:D:69:GLN:HB3	3:D:82:ASN:HB2	1.93	0.51
4:I:155:ARG:NH1	4:I:285:ILE:O	2.43	0.51
4:E:219:SER:HB2	4:E:224:SER:HB2	1.93	0.51
4:I:95:GLN:OE1	6:L:19:G:N1	2.39	0.50
4:F:35:LYS:HG3	4:F:118:LEU:HB2	1.92	0.50
4:J:122:GLY:HA3	4:J:140:VAL:HG11	1.93	0.50
4:J:207:ASP:HB2	4:J:210:LEU:HB3	1.94	0.50
4:F:185:GLN:O	4:G:236:ARG:NH1	2.44	0.50
4:I:32:PHE:HZ	4:I:136:LEU:HD21	1.75	0.50
4:E:43:MET:HG2	4:E:111:VAL:HG22	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:43:MET:CE	4:G:109:LEU:HD21	2.42	0.49
4:G:94:LEU:O	4:G:258:GLY:N	2.45	0.49
4:I:122:GLY:HA3	4:I:140:VAL:HG11	1.94	0.49
5:K:130:ARG:HD3	5:K:131:ILE:HG12	1.94	0.49
4:E:102:LEU:HD12	4:E:242:GLU:HA	1.94	0.49
4:F:63:ARG:NH2	4:F:101:ASN:OD1	2.46	0.49
5:K:63:SER:OG	5:K:64:ALA:N	2.46	0.49
3:D:200:LEU:HD22	3:D:210:LEU:HD12	1.94	0.49
4:G:249:LEU:HD13	4:H:66:SER:HB3	1.95	0.49
4:G:284:THR:HA	4:G:299:ALA:HA	1.93	0.49
4:H:284:THR:HA	4:H:299:ALA:HA	1.95	0.49
4:I:325:LEU:HD22	4:I:336:PRO:HB3	1.94	0.49
2:C:302:ARG:HD3	2:C:303:THR:HB	1.95	0.49
4:G:328:ASN:OD1	4:G:332:ARG:NH2	2.44	0.49
4:G:63:ARG:HH22	4:G:101:ASN:HD22	1.61	0.48
4:F:52:ASP:OD1	4:F:179:ARG:NH2	2.46	0.48
4:E:162:ASN:HB2	4:E:170:ARG:HH12	1.77	0.48
3:D:65:ARG:HB3	3:D:106:GLU:HB3	1.95	0.48
4:E:289:TYR:HB2	4:E:313:TYR:HE2	1.77	0.48
4:F:82:PRO:HA	4:F:85:LEU:HB3	1.96	0.48
1:A:39:CYS:SG	4:H:84:LYS:NZ	2.73	0.48
3:D:7:GLU:HG3	3:D:162:LEU:HD22	1.95	0.48
4:I:129:ASP:OD1	4:I:130:ALA:N	2.47	0.48
3:D:10:LEU:HB3	3:D:111:LEU:HB2	1.95	0.47
4:E:27:ALA:HB1	4:E:128:ASN:HB3	1.96	0.47
2:C:44:LEU:HA	2:C:48:ARG:HB3	1.95	0.47
4:H:277:LYS:NZ	6:L:22:U:OP1	2.40	0.47
4:I:63:ARG:NH2	4:I:101:ASN:OD1	2.46	0.47
4:F:41:ALA:HB1	4:F:111:VAL:HG13	1.96	0.47
4:H:298:ILE:HD13	4:H:313:TYR:HD2	1.79	0.47
4:I:170:ARG:NH2	4:I:196:ALA:O	2.47	0.47
3:D:282:ASN:ND2	6:L:3:A:H62	2.12	0.47
4:E:301:GLU:HB3	4:E:304:GLY:HA2	1.96	0.47
2:C:29:LEU:HD11	2:C:32:LEU:HD11	1.97	0.47
4:G:34:ARG:NH2	4:H:242:GLU:OE1	2.46	0.47
4:H:77:THR:HG22	4:H:80:ARG:HH22	1.79	0.47
3:D:64:HIS:NE2	3:D:108:SER:OG	2.34	0.47
5:K:39:ARG:HE	5:K:63:SER:HG	1.61	0.46
4:G:288:TRP:CD1	4:G:339:GLU:HB3	2.50	0.46
4:J:304:GLY:H	4:J:314:ARG:HB2	1.81	0.46
3:D:216:LEU:HB3	3:D:218:ARG:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:109:LEU:HD13	4:F:243:VAL:HG11	1.97	0.46
3:D:131:GLN:HB2	3:D:156:ASN:HD22	1.80	0.46
3:D:278:ARG:NH1	3:D:314:TYR:OH	2.48	0.46
4:G:51:ARG:HH11	4:G:52:ASP:H	1.62	0.46
4:E:338:VAL:HG13	4:E:341:GLN:HE21	1.81	0.46
2:C:338:ALA:HB1	3:D:182:PRO:HG3	1.98	0.46
4:G:226:HIS:CD2	4:H:174:GLU:HG2	2.51	0.46
4:H:328:ASN:O	4:H:334:GLU:N	2.45	0.46
3:D:2:SER:HA	3:D:320:ARG:HD3	1.97	0.46
4:J:305:SER:HA	4:J:312:ALA:HA	1.98	0.46
1:A:12:ASN:OD1	1:A:15:SER:OG	2.29	0.46
4:I:52:ASP:OD1	4:I:179:ARG:NH2	2.45	0.46
4:J:52:ASP:OD1	4:J:179:ARG:NH2	2.46	0.45
4:J:247:GLN:NE2	6:L:11:C:OP2	2.49	0.45
4:E:176:VAL:HB	4:E:196:ALA:HB3	1.98	0.45
2:C:72:PRO:HG3	2:C:251:ILE:HD11	1.98	0.45
2:C:329:ILE:HA	2:C:332:GLU:HG2	1.98	0.45
4:F:248:GLU:OE2	4:F:274:HIS:NE2	2.50	0.45
2:C:206:LEU:HD13	2:C:263:TRP:HH2	1.82	0.45
4:F:290:PRO:HB3	4:F:318:GLN:HE21	1.81	0.45
2:C:418:LEU:HA	2:C:421:GLU:HG3	1.98	0.45
4:H:54:SER:O	4:H:110:LYS:NZ	2.45	0.45
2:C:405:ILE:HG12	2:C:411:ALA:HB2	1.98	0.45
3:D:25:SER:HB3	3:D:252:PRO:HG3	1.99	0.45
1:A:2:LYS:HB3	1:A:2:LYS:HE2	1.46	0.45
4:I:245:PRO:HG2	4:I:264:LEU:HD22	1.99	0.44
5:K:106:LYS:O	6:L:56:G:N2	2.50	0.44
2:C:241:GLN:HB3	2:C:262:ASN:HB3	1.98	0.44
4:J:151:GLU:OE1	4:J:154:ARG:NH2	2.51	0.44
4:J:160:LEU:HD22	4:J:231:VAL:HG21	2.00	0.44
2:C:102:ASP:OD1	2:C:103:ARG:NH1	2.51	0.44
4:E:32:PHE:HZ	4:E:136:LEU:HD21	1.83	0.44
4:I:249:LEU:HD22	4:J:94:LEU:HB3	1.98	0.44
4:E:289:TYR:HB2	4:E:313:TYR:CE2	2.53	0.44
4:G:54:SER:HB2	4:G:110:LYS:HE3	2.00	0.44
4:E:327:ASP:OD2	4:F:74:ARG:NH1	2.45	0.43
4:G:44:SER:HB2	4:G:110:LYS:HB3	1.98	0.43
4:J:32:PHE:O	6:L:5:G:O2'	2.35	0.43
4:G:48:TRP:CE2	4:G:236:ARG:HD2	2.52	0.43
4:G:81:ASP:HB3	4:G:84:LYS:HB2	2.00	0.43
4:H:339:GLU:HA	4:H:342:HIS:CD2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:J:52:ASP:HA	4:J:179:ARG:HH12	1.83	0.43
4:G:77:THR:HA	4:G:80:ARG:HH12	1.82	0.43
4:H:164:ARG:HH21	4:H:201:LEU:HG	1.84	0.43
1:B:2:LYS:HB3	1:B:2:LYS:HE2	1.46	0.43
3:D:248:LEU:HB3	3:D:286:LEU:HD11	2.00	0.43
4:F:328:ASN:OD1	4:F:332:ARG:NH2	2.47	0.43
4:J:139:THR:HA	4:J:142:THR:HG22	2.00	0.43
4:E:102:LEU:HD23	4:E:102:LEU:HA	1.85	0.43
5:K:14:GLU:O	6:L:40:G:N2	2.52	0.43
1:B:12:ASN:OD1	1:B:15:SER:OG	2.29	0.43
2:C:362:TRP:CD1	2:C:363:LEU:HG	2.54	0.43
4:J:157:ALA:HB2	4:J:217:ILE:HD12	2.00	0.43
4:H:328:ASN:HB3	4:H:334:GLU:HB2	2.01	0.43
5:K:169:GLU:HA	5:K:179:SER:HB2	2.01	0.43
3:D:305:HIS:HB3	3:D:316:TRP:HB3	2.01	0.42
4:I:262:LYS:HE2	4:I:264:LEU:HD21	2.00	0.42
5:K:91:VAL:HA	5:K:92:PRO:HD3	1.94	0.42
5:K:95:THR:HA	5:K:185:PRO:HB3	2.01	0.42
2:C:285:SER:O	2:C:356:HIS:NE2	2.47	0.42
3:D:258:LEU:HD23	3:D:281:GLU:HB3	2.01	0.42
4:I:123:THR:HA	4:I:137:LEU:HD21	1.99	0.42
4:F:77:THR:HG22	4:F:80:ARG:HH12	1.84	0.42
4:G:207:ASP:HB3	4:G:210:LEU:HB2	2.00	0.42
2:C:241:GLN:HG3	3:D:99:GLU:OE2	2.20	0.42
4:G:56:GLU:O	4:G:56:GLU:CG	2.66	0.42
2:C:335:GLN:NE2	3:D:181:LEU:O	2.53	0.42
4:E:52:ASP:OD1	4:E:179:ARG:NH1	2.47	0.42
4:I:179:ARG:NH2	4:I:193:ARG:HH12	2.17	0.42
4:E:102:LEU:HA	4:E:103:PRO:HD3	1.82	0.42
4:G:268:ARG:HH22	4:H:103:PRO:HB3	1.84	0.42
5:K:2:ASP:O	5:K:4:TYR:N	2.52	0.42
3:D:272:ASP:OD1	3:D:272:ASP:N	2.53	0.42
4:F:151:GLU:HG2	4:F:154:ARG:HH21	1.85	0.42
4:H:323:TYR:OH	4:I:72:SER:O	2.31	0.42
4:I:129:ASP:HB3	4:I:132:TYR:HB3	2.01	0.42
3:D:167:GLU:OE1	3:D:170:ARG:NH2	2.53	0.41
4:F:28:SER:OG	4:F:128:ASN:ND2	2.52	0.41
2:C:321:ARG:HH11	2:C:322:ARG:HG3	1.85	0.41
4:E:320:LEU:HD13	4:E:325:LEU:HD11	2.02	0.41
4:G:29:VAL:HB	4:G:128:ASN:HB2	2.01	0.41
2:C:66:VAL:HG11	2:C:81:ASN:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:197:LEU:HD12	3:D:210:LEU:HD13	2.02	0.41
4:E:35:LYS:HG3	4:E:118:LEU:HB2	2.02	0.41
4:F:339:GLU:HA	4:F:342:HIS:HD2	1.85	0.41
4:H:77:THR:HA	4:H:80:ARG:HH12	1.84	0.41
4:J:35:LYS:HG3	4:J:118:LEU:HB2	2.01	0.41
4:J:284:THR:HA	4:J:299:ALA:HA	2.03	0.41
1:B:16:GLU:HG2	4:J:257:LYS:HE3	2.03	0.41
4:E:162:ASN:HD22	4:E:201:LEU:HD12	1.86	0.41
4:H:304:GLY:H	4:H:314:ARG:HB3	1.85	0.41
4:E:141:ALA:HA	4:E:144:VAL:HG12	2.02	0.41
4:H:317:LYS:HZ2	4:I:82:PRO:HD3	1.86	0.41
4:I:321:ASP:N	4:I:321:ASP:OD1	2.51	0.41
1:B:2:LYS:NZ	1:B:23:GLU:HB3	2.36	0.41
2:C:272:TRP:HH2	3:D:70:ILE:HD11	1.85	0.41
2:C:375:LEU:HB2	2:C:377:ARG:HG2	2.03	0.41
2:C:386:GLU:O	2:C:390:ARG:HD3	2.21	0.41
4:E:150:ALA:HA	4:E:153:ALA:HB3	2.02	0.41
4:F:259:GLN:HE21	4:F:263:THR:HA	1.86	0.41
4:I:43:MET:HB3	4:I:109:LEU:HD11	2.03	0.41
4:G:312:ALA:N	4:H:86:ASP:OD1	2.49	0.40
4:J:182:HIS:CD2	4:J:216:LEU:HD21	2.56	0.40
4:G:43:MET:HE2	4:G:109:LEU:HD21	2.03	0.40
4:H:68:ARG:NE	6:L:26:G:O2'	2.54	0.40
4:I:288:TRP:CD1	4:I:339:GLU:HB3	2.56	0.40
1:A:2:LYS:NZ	1:A:23:GLU:HB3	2.36	0.40
2:C:268:LEU:HD13	2:C:272:TRP:CD2	2.56	0.40
4:F:154:ARG:O	4:F:158:HIS:ND1	2.55	0.40
4:E:54:SER:O	4:E:110:LYS:NZ	2.49	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	66/68 (97%)	62 (94%)	4 (6%)	0	100	100
1	B	66/68 (97%)	62 (94%)	4 (6%)	0	100	100
2	C	420/434 (97%)	368 (88%)	51 (12%)	1 (0%)	47	69
3	D	302/327 (92%)	281 (93%)	21 (7%)	0	100	100
4	E	279/360 (78%)	248 (89%)	31 (11%)	0	100	100
4	F	324/360 (90%)	298 (92%)	26 (8%)	0	100	100
4	G	323/360 (90%)	299 (93%)	24 (7%)	0	100	100
4	H	333/360 (92%)	313 (94%)	20 (6%)	0	100	100
4	I	323/360 (90%)	303 (94%)	20 (6%)	0	100	100
4	J	335/360 (93%)	314 (94%)	20 (6%)	1 (0%)	41	62
5	K	185/187 (99%)	151 (82%)	34 (18%)	0	100	100
All	All	2956/3244 (91%)	2699 (91%)	255 (9%)	2 (0%)	54	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	J	333	ASP
2	C	92	PRO

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	58/58 (100%)	52 (90%)	6 (10%)	7	12
1	B	58/58 (100%)	52 (90%)	6 (10%)	7	12
2	C	353/365 (97%)	345 (98%)	8 (2%)	50	73
3	D	255/270 (94%)	252 (99%)	3 (1%)	71	86
4	E	221/290 (76%)	218 (99%)	3 (1%)	67	84
4	F	261/290 (90%)	260 (100%)	1 (0%)	91	97
4	G	260/290 (90%)	254 (98%)	6 (2%)	50	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	H	266/290 (92%)	263 (99%)	3 (1%)	73	88
4	I	260/290 (90%)	256 (98%)	4 (2%)	65	82
4	J	268/290 (92%)	266 (99%)	2 (1%)	84	93
5	K	160/160 (100%)	156 (98%)	4 (2%)	47	70
All	All	2420/2651 (91%)	2374 (98%)	46 (2%)	59	77

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	18	GLU
1	A	23	GLU
1	A	36	LYS
1	A	53	LEU
1	A	63	ARG
1	B	2	LYS
1	B	18	GLU
1	B	23	GLU
1	B	36	LYS
1	B	53	LEU
1	B	63	ARG
2	C	222	ARG
2	C	256	ASN
2	C	293	ARG
2	C	302	ARG
2	C	319	ARG
2	C	321	ARG
2	C	341	ARG
2	C	424	MET
3	D	28	THR
3	D	90	ARG
3	D	218	ARG
4	E	168	ARG
4	E	170	ARG
4	E	190	ARG
4	F	280	ASN
4	G	44	SER
4	G	50	GLN
4	G	51	ARG
4	G	55	GLN
4	G	56	GLU

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Mol	Chain	Res	Type
4	G	57	TRP
4	H	68	ARG
4	H	145	ASN
4	H	317	LYS
4	I	128	ASN
4	I	205	LYS
4	I	280	ASN
4	I	283	ARG
4	J	78	LYS
4	J	137	LEU
5	K	57	ARG
5	K	73	ARG
5	K	130	ARG
5	K	156	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	237	ASN
2	C	256	ASN
2	C	283	HIS
2	C	369	GLN
3	D	282	ASN
3	D	293	HIS
4	E	128	ASN
4	E	328	ASN
4	E	341	GLN
4	F	128	ASN
4	F	259	GLN
4	F	318	GLN
4	G	50	GLN
4	G	90	GLN
4	G	101	ASN
4	G	182	HIS
4	G	226	HIS
4	H	128	ASN
4	H	145	ASN
4	H	318	GLN
4	H	328	ASN
4	I	128	ASN
4	I	280	ASN
4	J	90	GLN

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Mol	Chain	Res	Type
4	J	147	GLN
4	J	241	GLN
5	K	29	HIS
5	K	154	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
6	L	59/60 (98%)	33 (55%)	1 (1%)

All (33) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
6	L	3	A
6	L	9	U
6	L	14	G
6	L	15	G
6	L	16	C
6	L	20	C
6	L	21	U
6	L	22	U
6	L	26	G
6	L	27	U
6	L	28	C
6	L	29	C
6	L	30	G
6	L	32	G
6	L	33	U
6	L	34	C
6	L	38	C
6	L	39	U
6	L	40	G
6	L	41	G
6	L	44	C
6	L	45	A
6	L	47	U
6	L	48	G
6	L	49	C
6	L	50	C
6	L	51	G
6	L	52	U

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Mol	Chain	Res	Type
6	L	53	A
6	L	57	G
6	L	58	C
6	L	59	A
6	L	60	G

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
6	L	58	C

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.

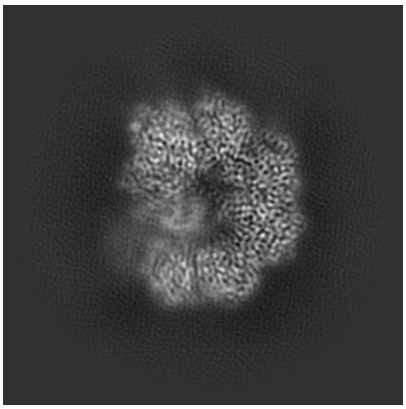
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-21358. These allow visual inspection of the internal detail of the map and identification of artifacts.

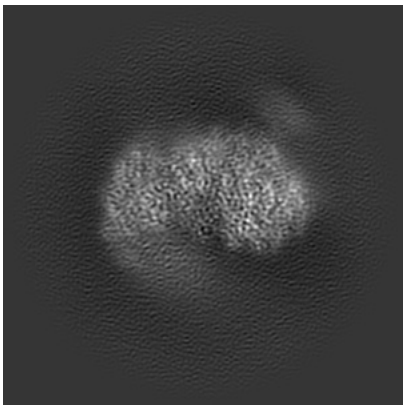
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

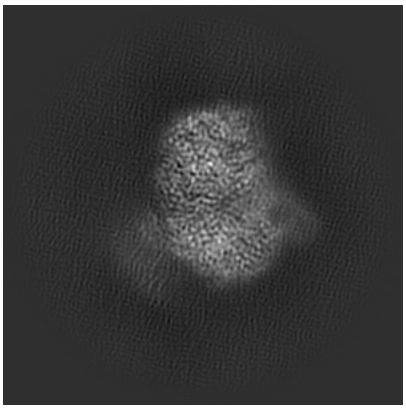
6.1.1 Primary map



X

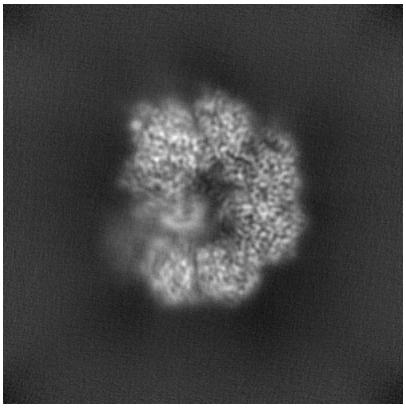


Y

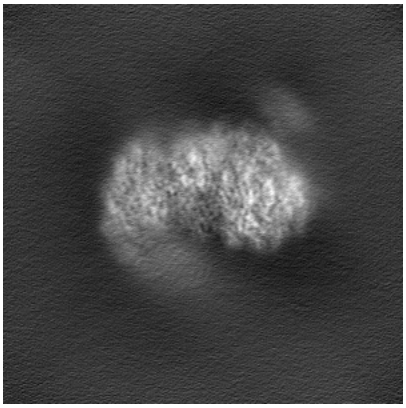


Z

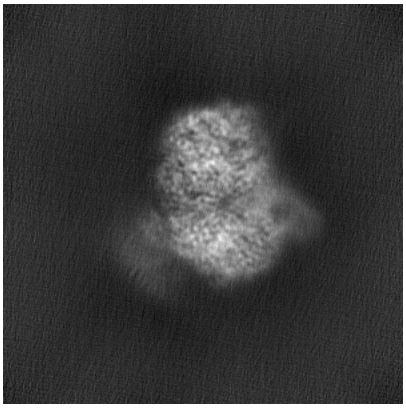
6.1.2 Raw map



X



Y

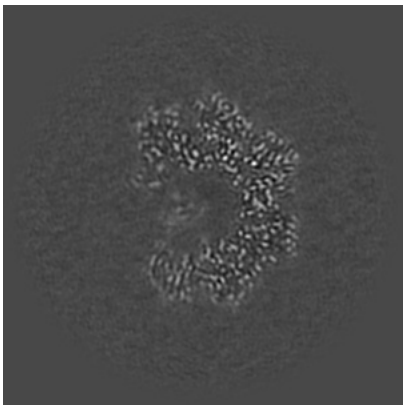


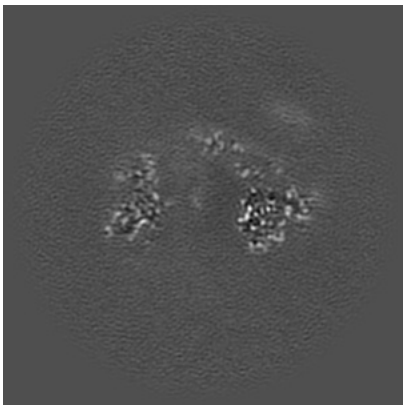
Z

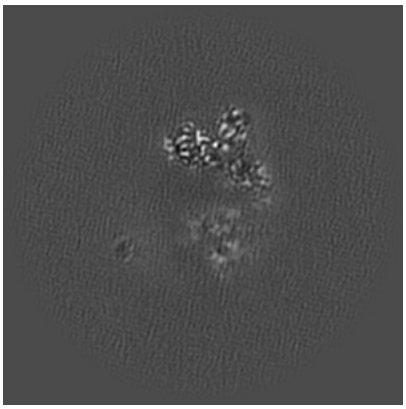
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

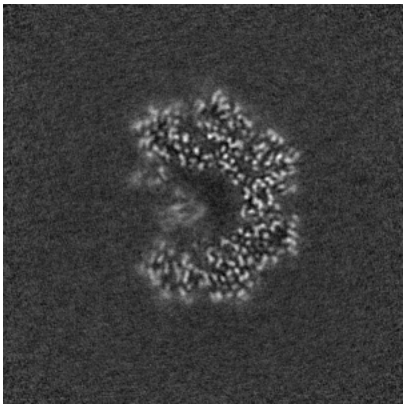
6.2.1 Primary map

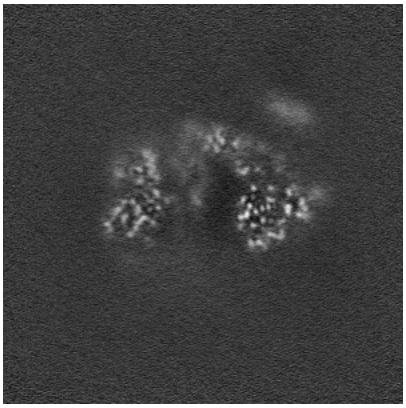

X Index: 192

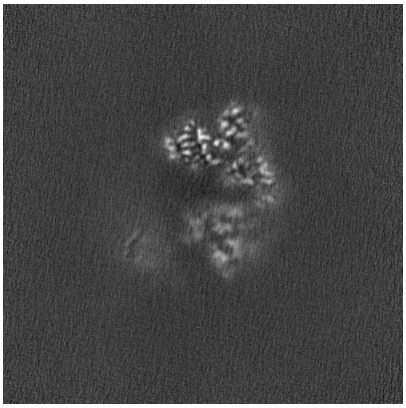

Y Index: 192


Z Index: 192

6.2.2 Raw map


X Index: 192

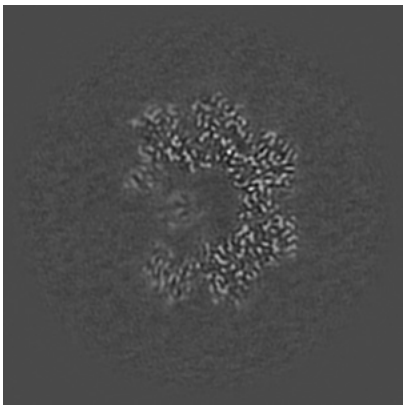

Y Index: 192

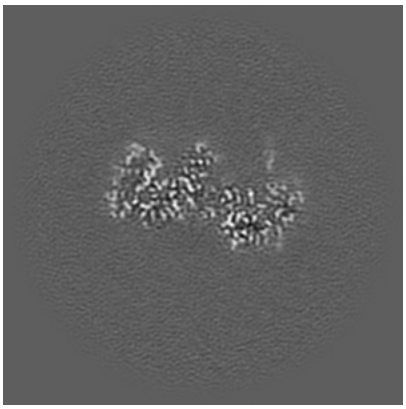

Z Index: 192

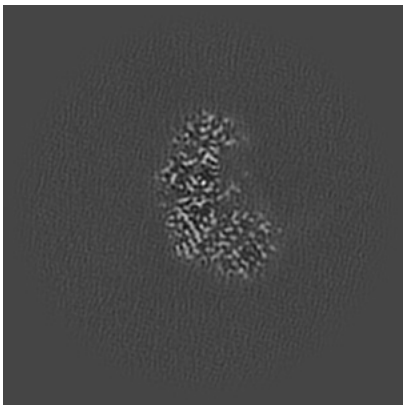
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

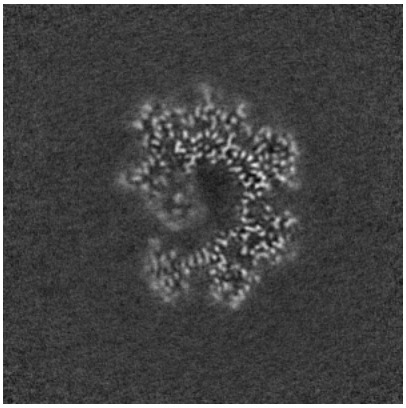
6.3.1 Primary map

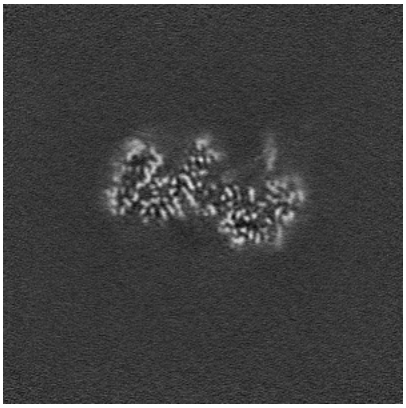

X Index: 194

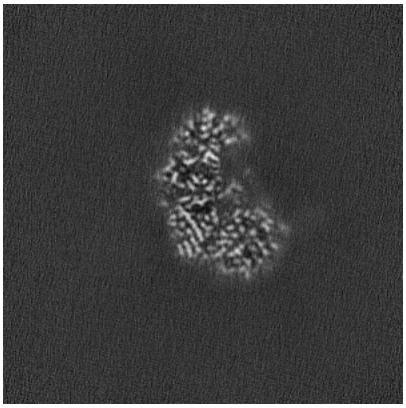

Y Index: 228


Z Index: 236

6.3.2 Raw map


X Index: 201

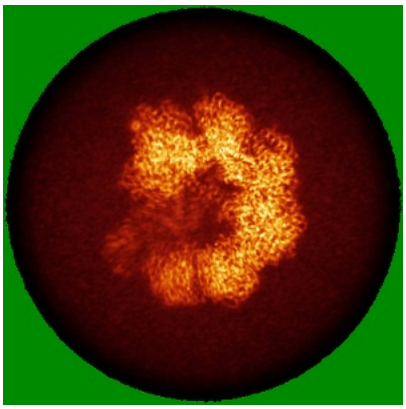

Y Index: 229

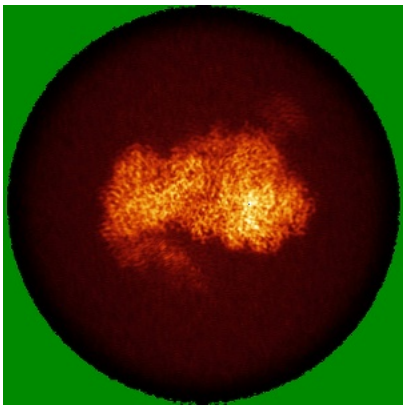

Z Index: 236

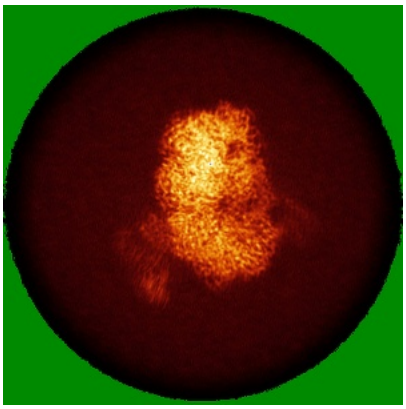
The images above show the largest variance slices of the map in three orthogonal directions.

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

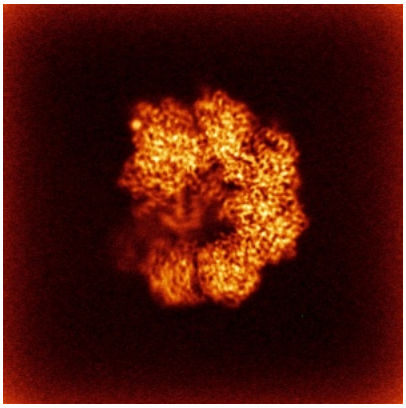
6.4.1 Primary map

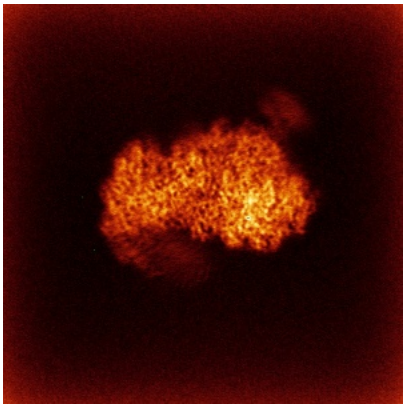

X

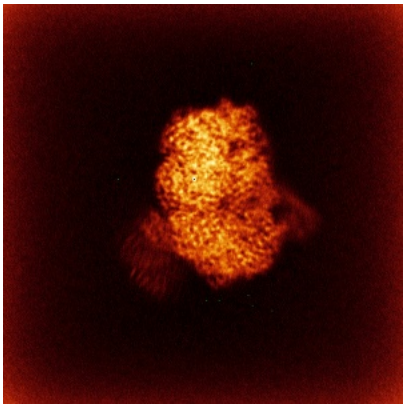

Y


Z

6.4.2 Raw map


X

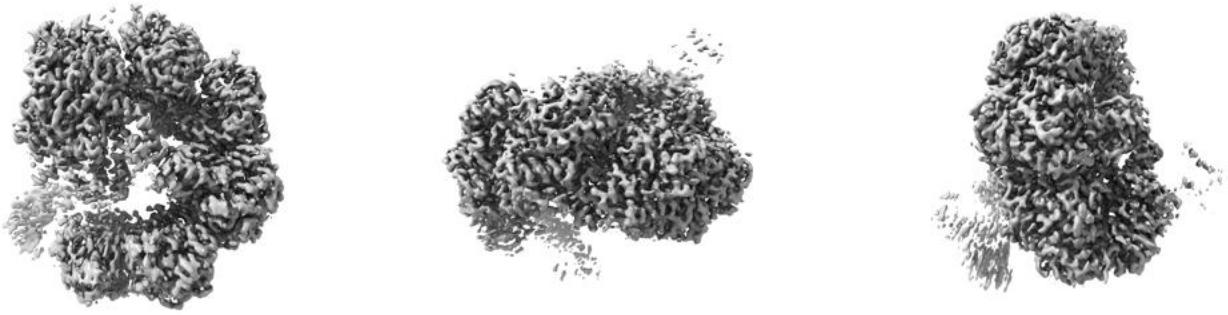

Y


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

6.5.2 Raw map

X Y Z

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

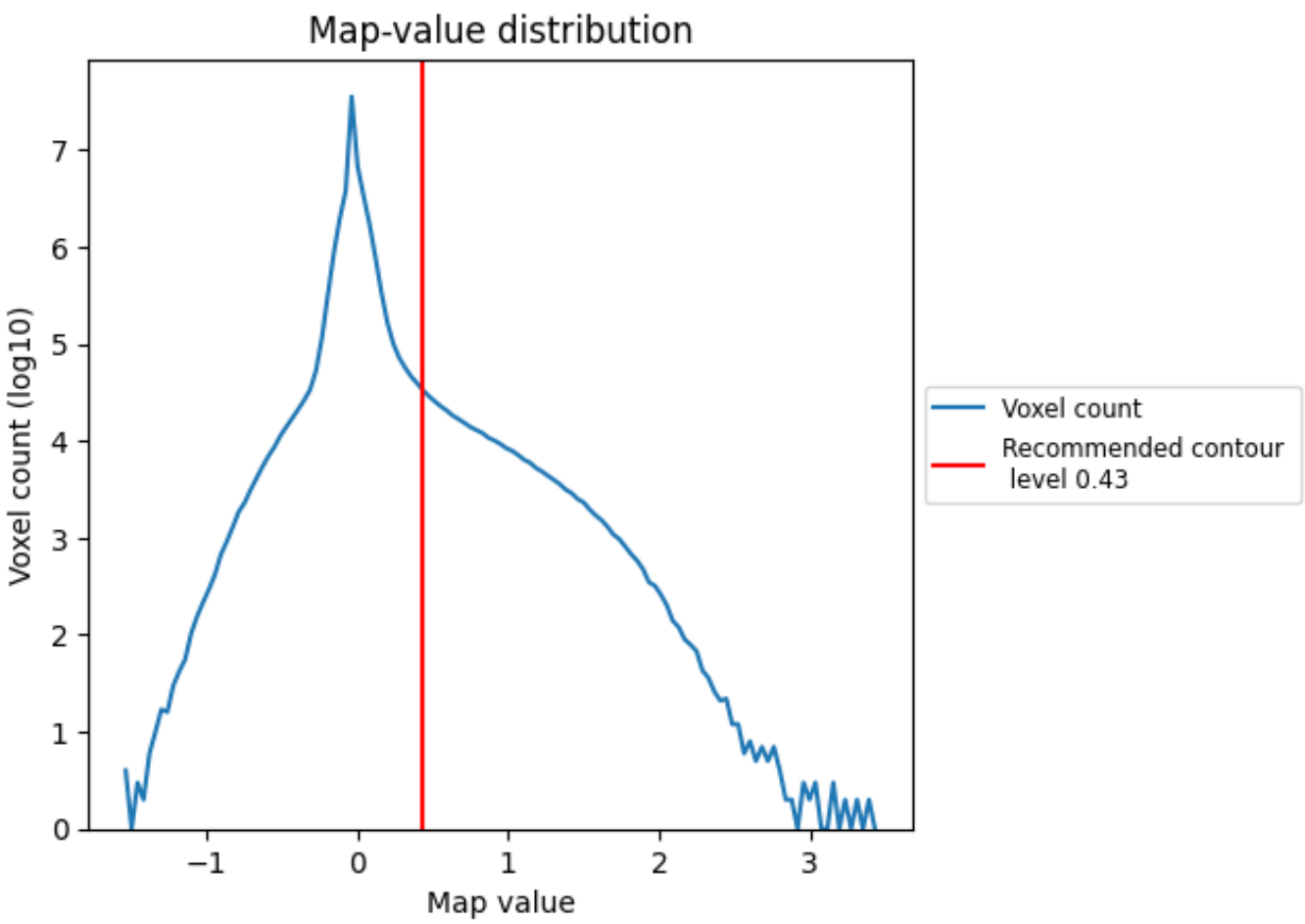
6.6 Mask visualisation (i)

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

7 Map analysis (i)

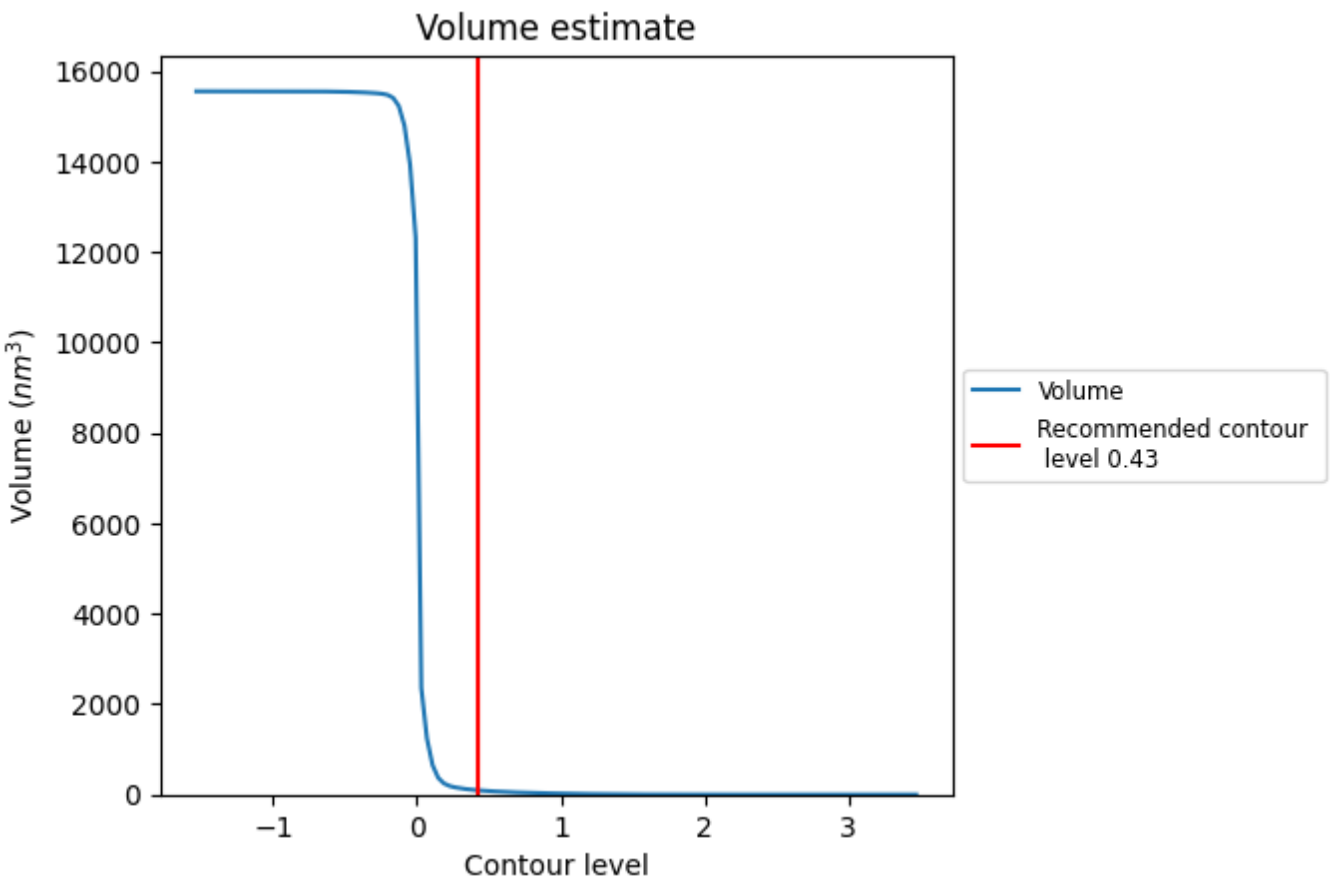
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution (i)



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

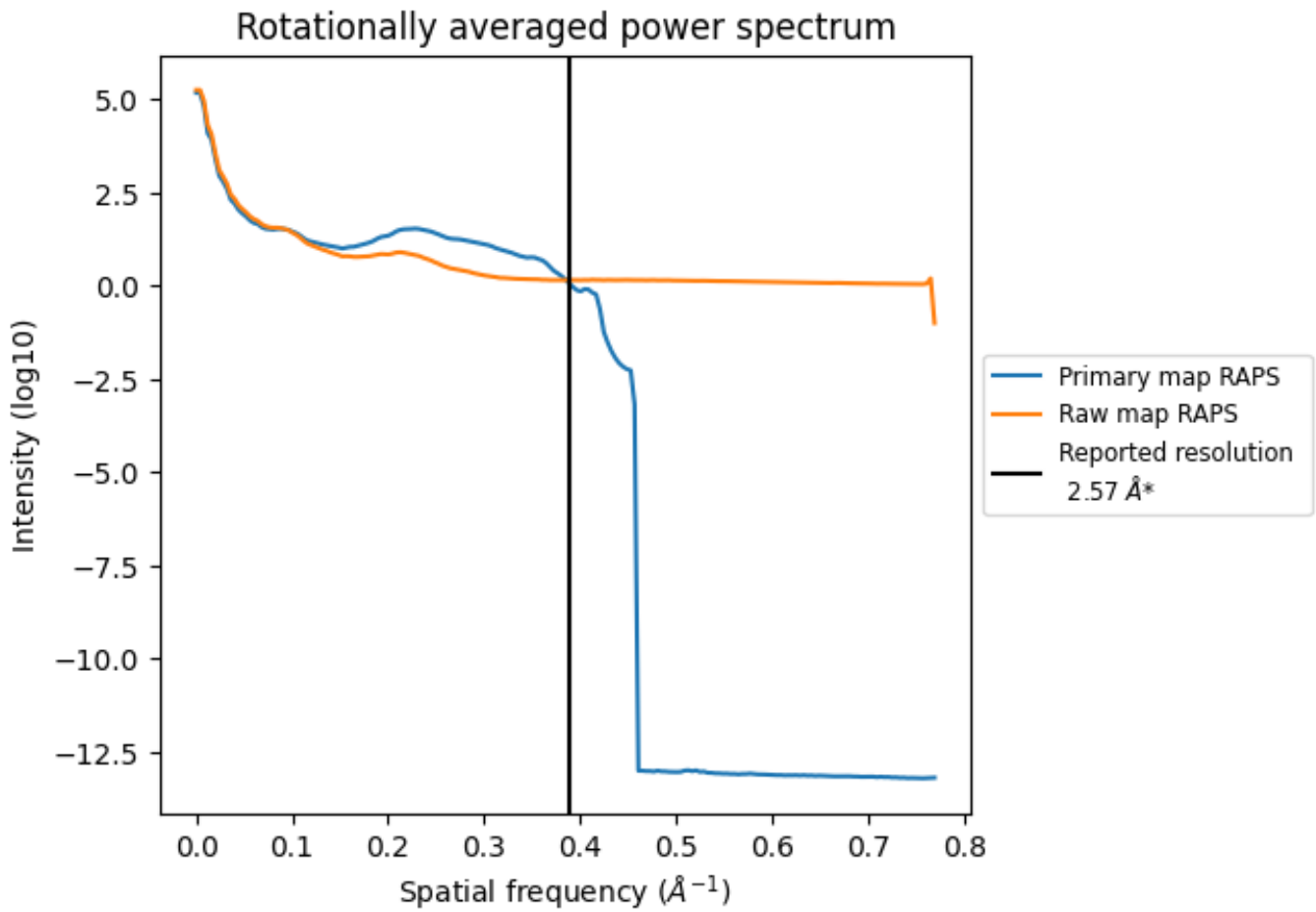
7.2 Volume estimate (i)



The volume at the recommended contour level is 93 nm^3; this corresponds to an approximate mass of 84 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)

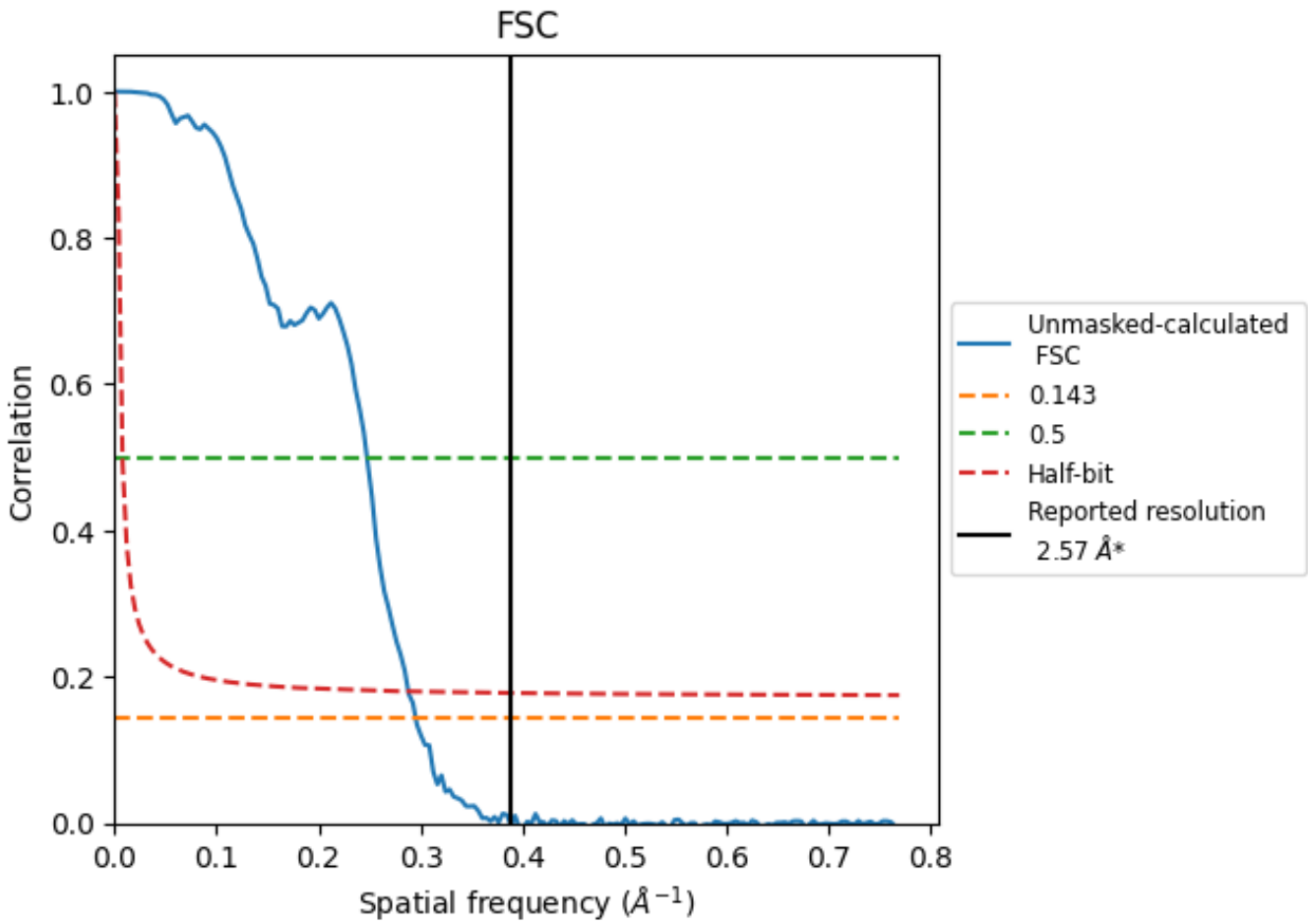


*Reported resolution corresponds to spatial frequency of 0.389 Å^{-1}

8 Fourier-Shell correlation (i)

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

8.1 FSC (i)



*Reported resolution corresponds to spatial frequency of 0.389 Å^{-1}

8.2 Resolution estimates (i)

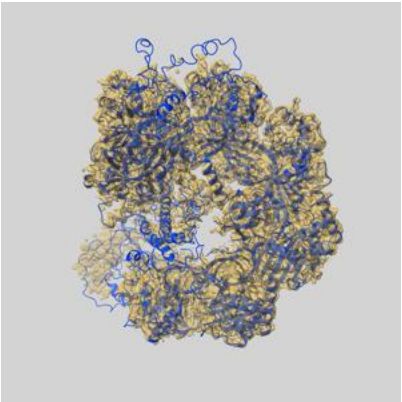
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.57	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.39	4.04	3.47

*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 3.39 differs from the reported value 2.57 by more than 10 %

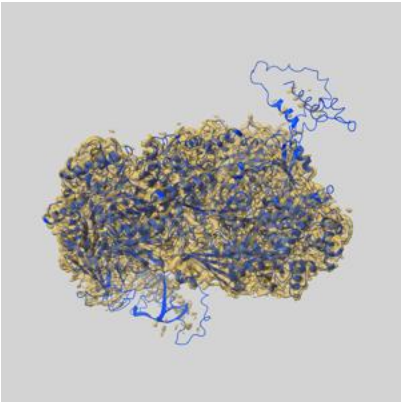
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-21358 and PDB model 6VQV. Per-residue inclusion information can be found in section 3 on page 9.

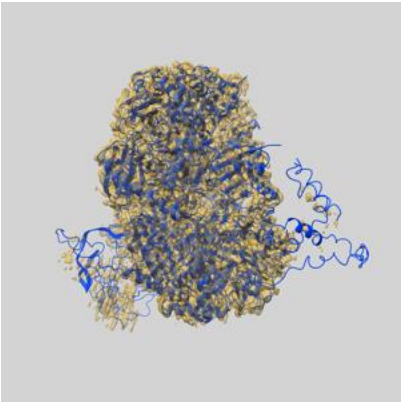
9.1 Map-model overlay (i)



X



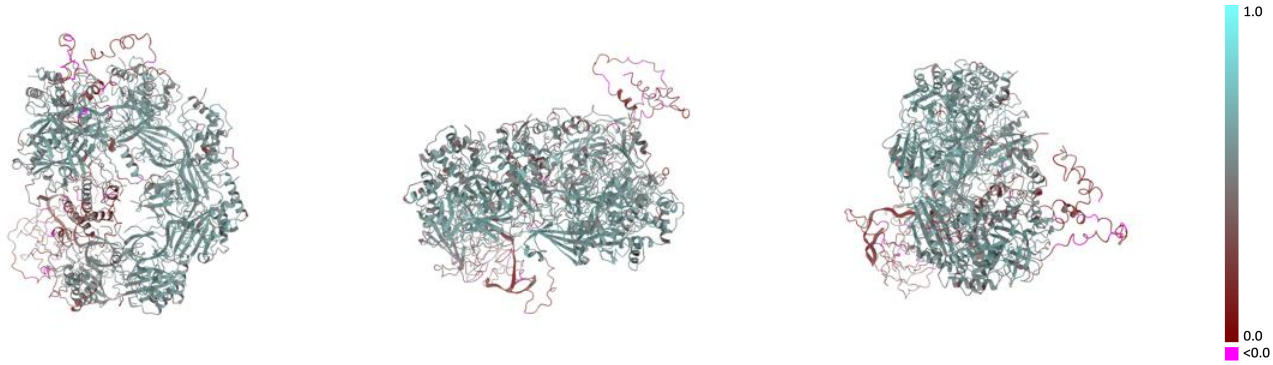
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.43 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.

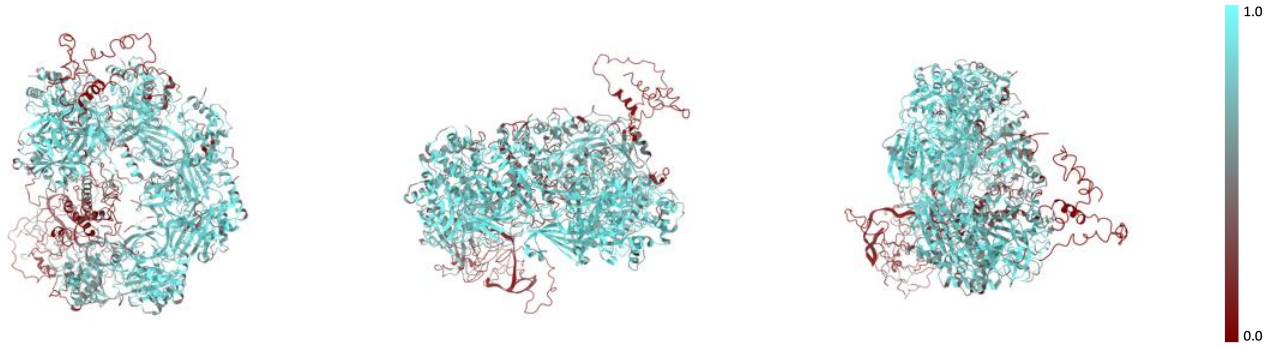
9.2 Q-score mapped to coordinate model (i)



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.

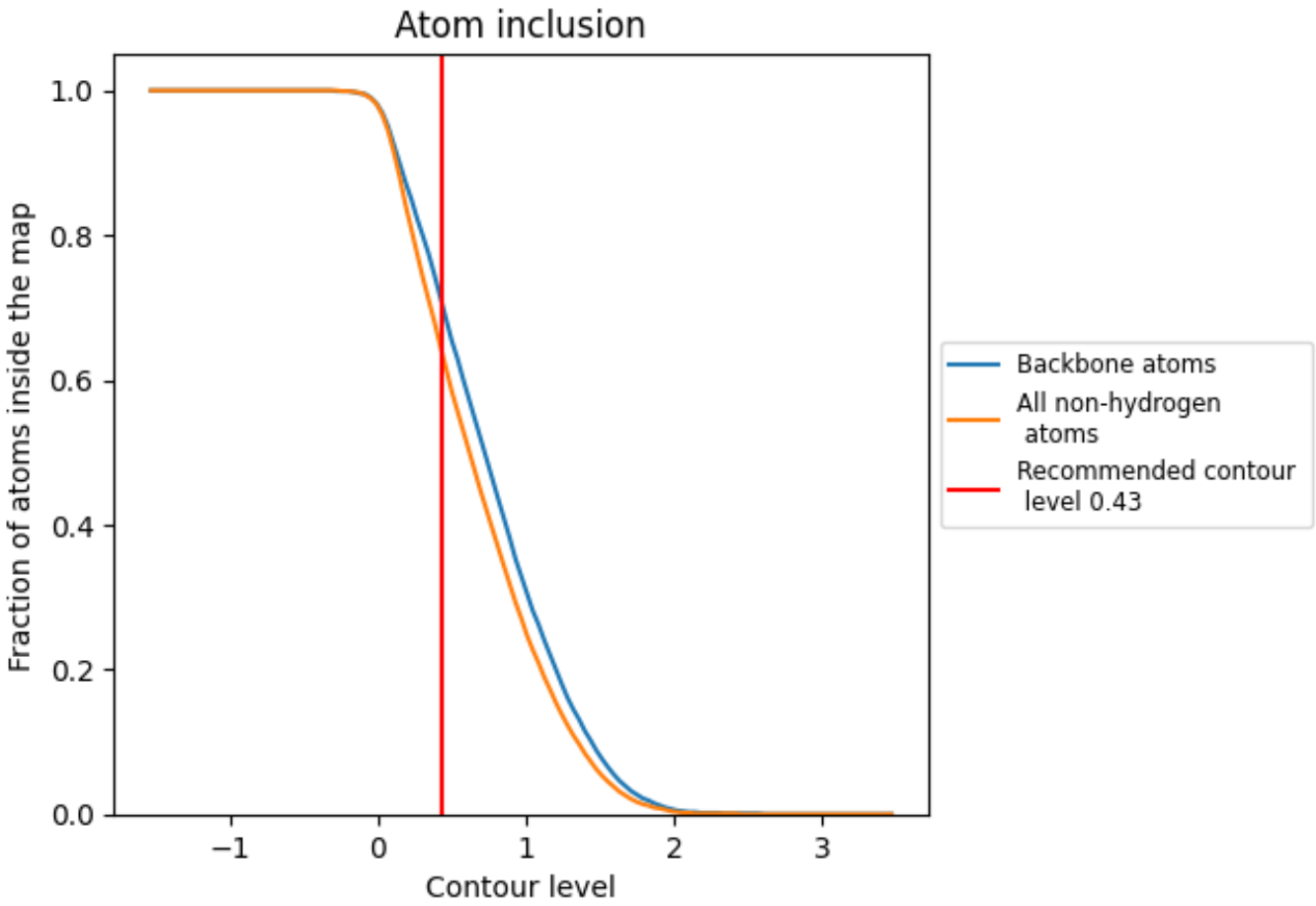
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.43).

9.4 Atom inclusion (i)



At the recommended contour level, 71% of all backbone atoms, 64% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.43) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6400	0.4950
A	0.7190	0.5370
B	0.6020	0.5080
C	0.3120	0.3430
D	0.7020	0.5380
E	0.5570	0.4510
F	0.7580	0.5460
G	0.8220	0.5830
H	0.8100	0.5700
I	0.8270	0.5850
J	0.7690	0.5630
K	0.0750	0.2140
L	0.6080	0.4540