



wwPDB EM Validation Summary Report ⓘ

Mar 9, 2026 – 04:29 AM UTC

PDB ID : 7XG4 / pdb_00007xg4
EMDB ID : EMD-33185
Title : CryoEM structure of type IV-A CasDinG bound NTS-nicked Csf-crRNA-dsDNA quaternary complex in a second state
Authors : Zhang, J.T.; Cui, N.; Huang, H.D.; Jia, N.
Deposited on : 2022-04-02
Resolution : 3.70 Å(reported)

This is a wwPDB EM Validation Summary 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:

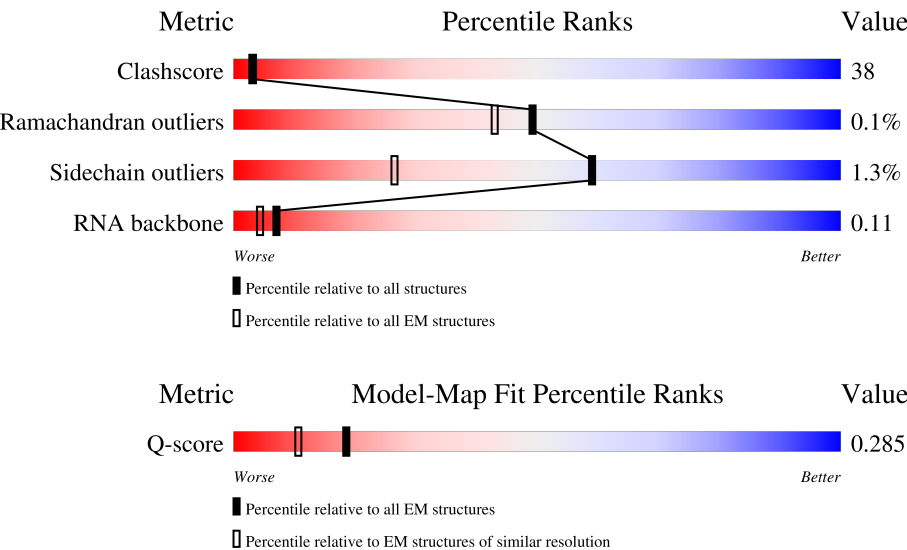
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 i

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

The reported resolution of this entry is 3.70 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.



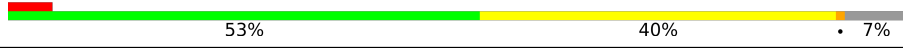
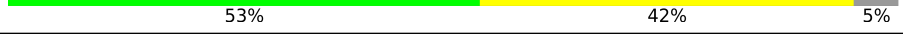
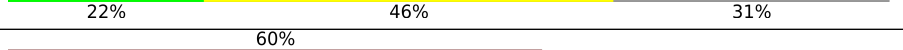
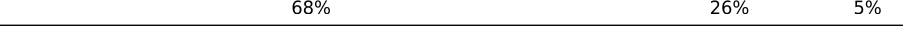
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	11569 (3.20 - 4.20)

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	253	6% 52% 42% 5%
2	B	220	5% 58% 40%
3	C	348	62% 32% 5%

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Mol	Chain	Length	Quality of chain
3	D	348	
3	E	348	
3	F	348	
3	G	348	
4	H	268	
5	I	61	
6	J	38	
7	K	54	
8	L	626	

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
5	23G	I	61	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 24769 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 Csf1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	241	Total	C	N	O	S	0	0
			1897	1202	336	347	12		

- Molecule 2 is a protein called Csf3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	219	Total	C	N	O	S	0	0
			1709	1080	316	305	8		

- Molecule 3 is a protein called Csf2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	329	Total	C	N	O	S	0	0
			2488	1570	439	467	12		
3	D	331	Total	C	N	O	S	0	0
			2503	1580	441	470	12		
3	E	324	Total	C	N	O	S	0	0
			2450	1546	431	461	12		
3	F	324	Total	C	N	O	S	0	0
			2450	1546	431	461	12		
3	G	280	Total	C	N	O	S	0	0
			2131	1352	370	398	11		

- Molecule 4 is a protein called Csf5.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	234	Total	C	N	O	S	0	0
			1800	1149	322	322	7		

- Molecule 5 is a RNA chain called crRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	I	61	Total	C	N	O	P	0	0
			1317	586	247	423	61		

- Molecule 6 is a DNA chain called NTS.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	J	36	Total	C	N	O	P	0	0
			722	355	98	233	36		

- Molecule 7 is a DNA chain called TS.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	K	37	Total	C	N	O	P	0	0
			759	361	137	224	37		

- Molecule 8 is a protein called Csf4.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	L	594	Total	C	N	O	S	0	0
			4542	2872	817	834	19		

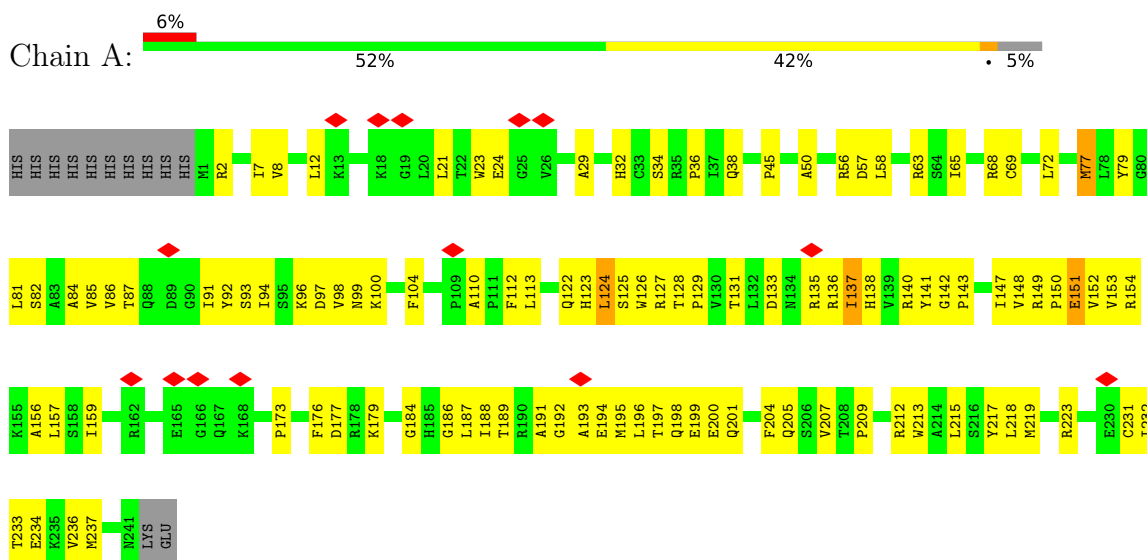
- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
9	A	1	Total	Zn	0
			1	1	

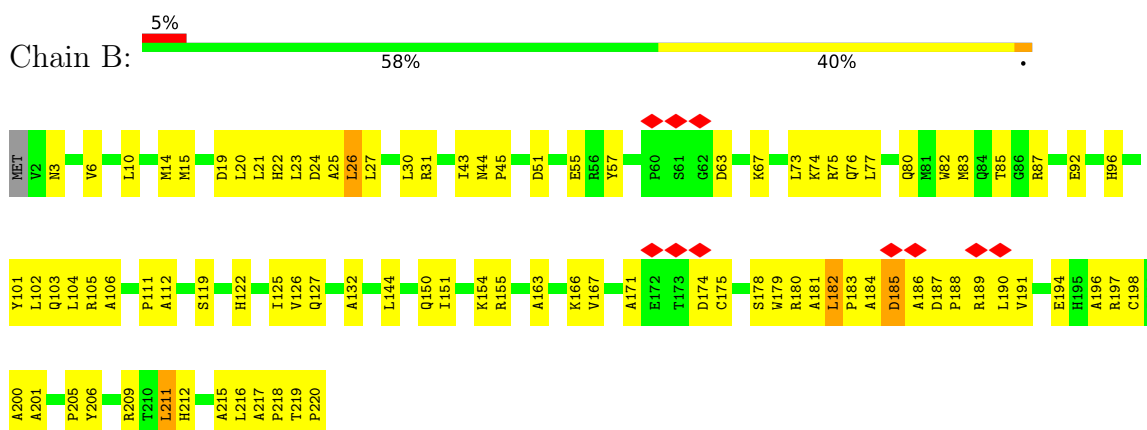
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Csf1

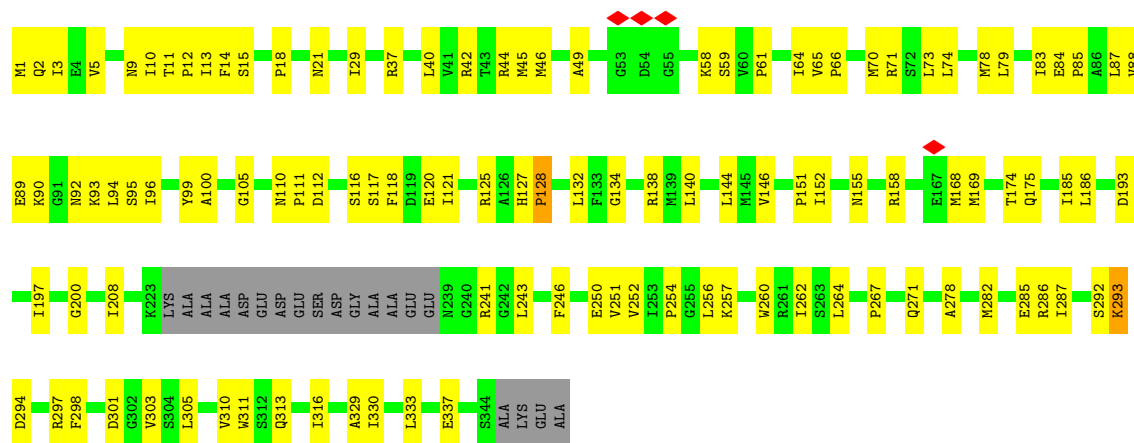


• Molecule 2: Csf3



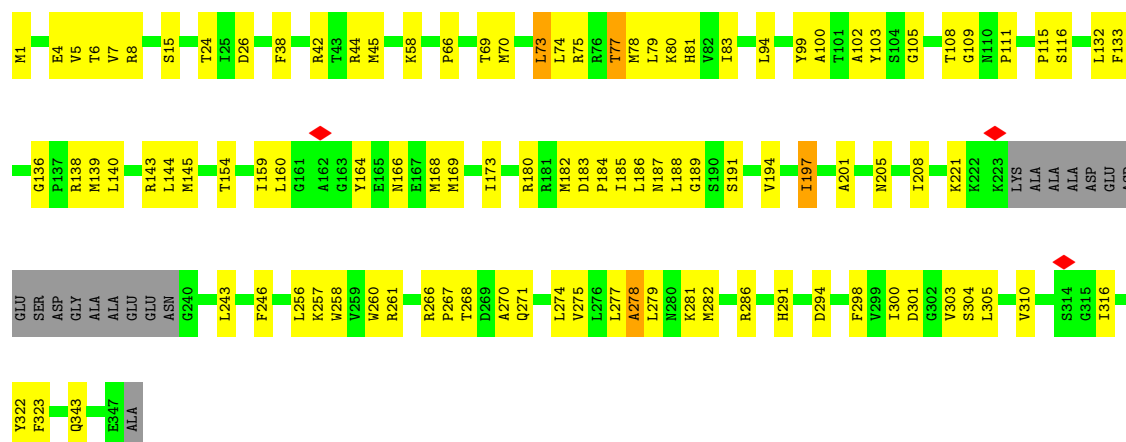
• Molecule 3: Csf2





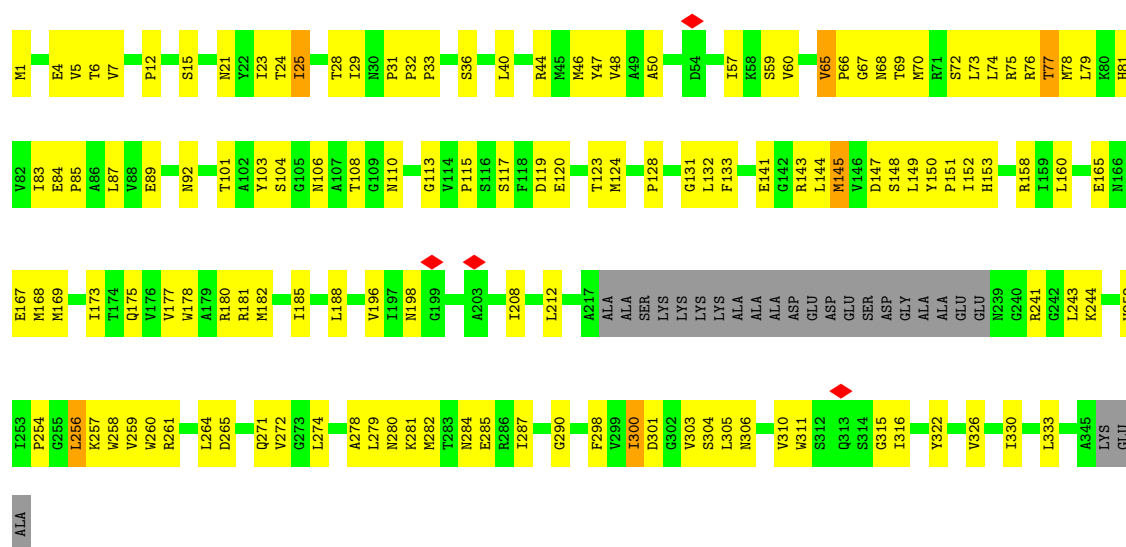
• Molecule 3: Csf2

Chain D: 66% 28% 5%

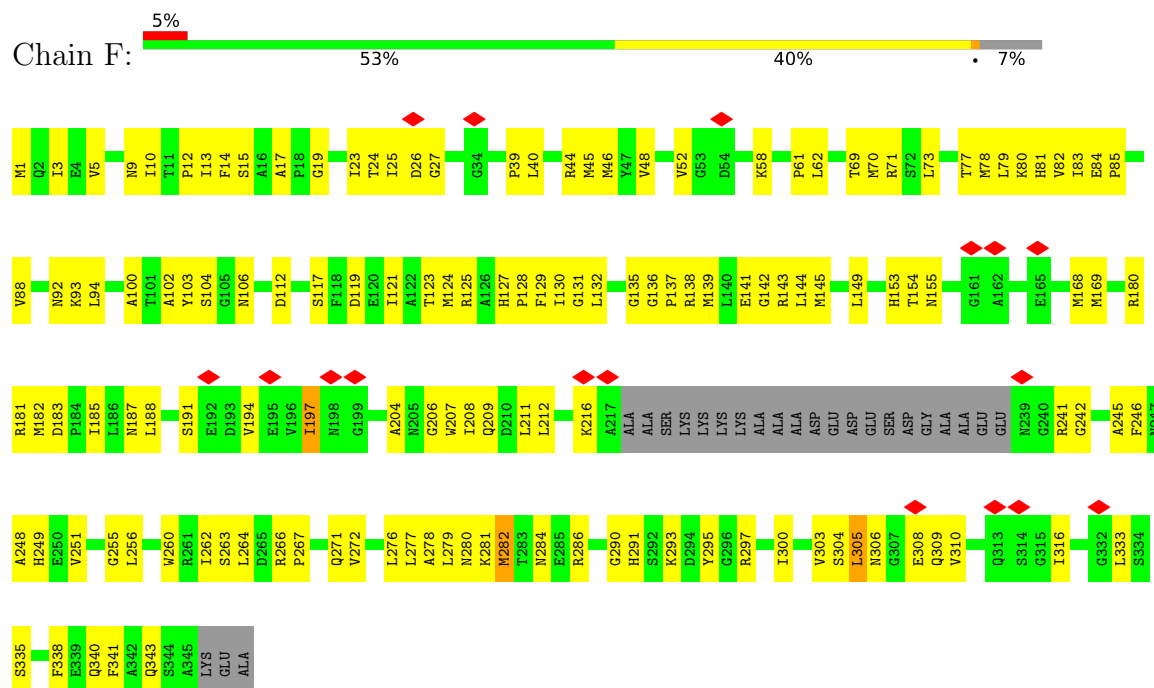


• Molecule 3: Csf2

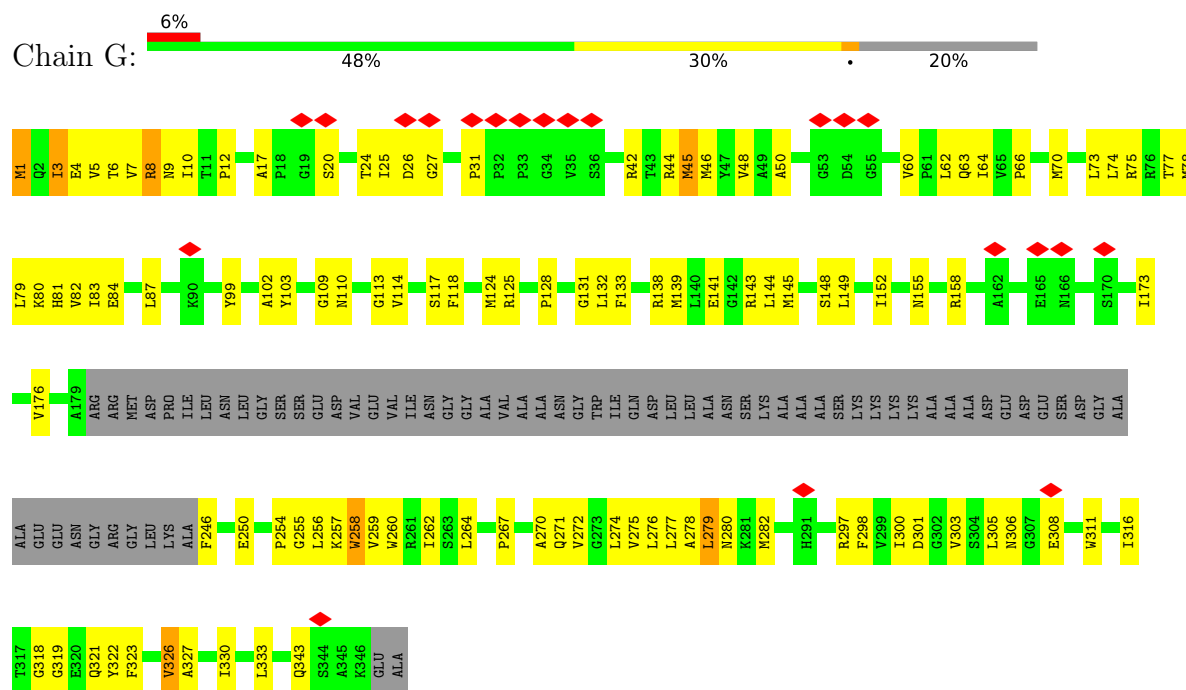
Chain E: 55% 37% 7%



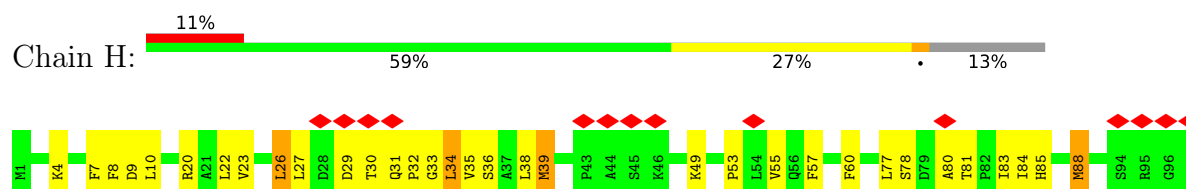
- Molecule 3: Csf2

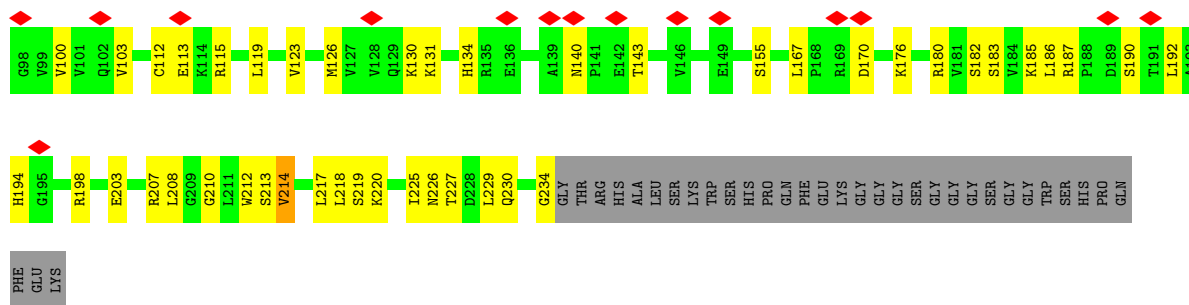


- Molecule 3: Csf2



- Molecule 4: Csf5





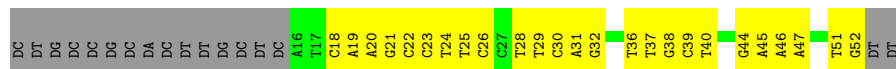
• Molecule 5: crRNA



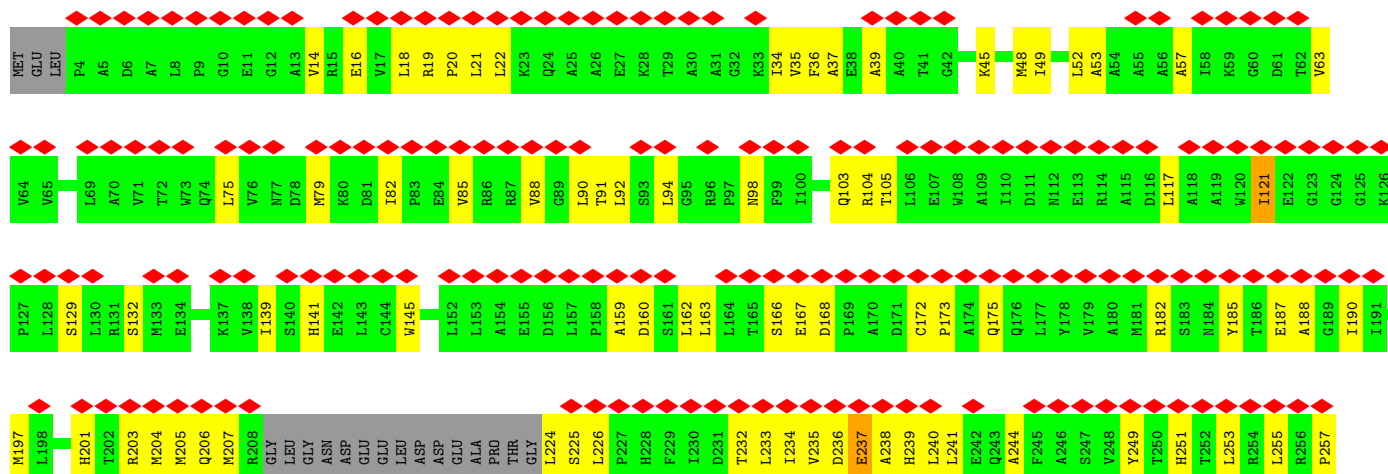
• Molecule 6: NTS



• Molecule 7: TS



• Molecule 8: Csf4



L599	R605	I606	Y607	G608	G609	A610	A611	W612	V613	A614	P615	P616	R617	Q618	I619	L620	D621	R622	Y623	K624	W625	ALA																																					
R513	A522	G525	H531	S532	L533	P534	D535	N536	P537	E538	L539	D540	R541	S544	D545	L546	V547	I548	T549	R550	Q555	N556	R557	S558	L559	T560	H561	A566	R571	I572	I573	S574	Q575	W579	G584	L585	L588	V589	R590	G593	H596	K597	N598																
A445	T446	E447	C448	A449	Q450	T451	T452	Q453	G454	V455	A456	S457	T458	A459	T463	L464	C467	T468	S469	Y470	Q471	N472	T473	E474	L475	L476	A477	G478	R479	L480	G481	A482	A483	L484	G485	D486	R487	L488	Q491	S492	K493	T494	S495	S496	A497	A498	T499	C500	A506	K507	H508	K509	A510	G511	I512				
W378	N379	A380	T381	G382	G383	A384	T385	L386	V387	S388	L391	F392	T393	T394	G395	D396	N397	G398	S399	L400	T401	K404	L405	E406	V407	P408	T409	E410	R411	F414	L415	P416	P417	V418	A421	A425	L428	L429	H430	K431	E432	F433	C434	A435	H436	E437	P438	D439	D440	S441	P442	E443	W444						
G318	A319	L320	P321	T322	K323	G324	M325	S326	R327	D328	A329	R330	S331	V332	I333	R334	I335	A336	T337	R338	A339	A340	N341	D342	A343	L344	S345	G346	H347	S348	R349	L350	R351	I352	E353	V354	T355	P356	V357	H358	S359	Y360	P361	M362	L363	L364	S365	G366	R367	S368	N369	L370	Q371	R372	A373	L374	L375	G376	L377
L258	M259	R260	T261	T262	E263	G264	LEU	GLY	SER	ARG	GLY	ARG	LYS	PRO	ALA	L274	D275	A276	L277	K278	E279	L280	F281	T282	Q283	M284	Q285	V286	A287	SER	ALA	ARG	SER	T292	N293	T294	S295	L296	N297	V298	P299	L300	S301	D302	V303	P304	E305	L306	I307	P308	A309	L310	K311	D312	T313	V314	K315	T316	L317

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	8882	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.072	Depositor
Minimum map value	-0.031	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.008	Depositor
Map size ($\text{\AA}$)	396.0, 396.0, 396.0	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.1, 1.1, 1.1	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 23G, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.61	1/1945 (0.1%)	0.86	0/2643
2	B	0.54	1/1754 (0.1%)	0.71	0/2384
3	C	0.48	0/2537	0.75	1/3439 (0.0%)
3	D	0.49	2/2552 (0.1%)	0.74	0/3458
3	E	0.53	1/2499 (0.0%)	0.80	2/3391 (0.1%)
3	F	0.52	0/2499	0.78	1/3391 (0.0%)
3	G	0.59	0/2177	0.91	0/2955
4	H	0.50	2/1835 (0.1%)	0.72	1/2475 (0.0%)
5	I	0.32	0/1446	0.78	4/2256 (0.2%)
6	J	0.30	0/800	0.72	0/1231
7	K	0.33	0/850	0.66	0/1310
8	L	0.47	3/4632 (0.1%)	0.68	2/6301 (0.0%)
All	All	0.50	10/25526 (0.0%)	0.76	11/35234 (0.0%)

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	151	ILE	CG1-CD1	-8.52	1.18	1.51
8	L	205	MET	SD-CE	-7.41	1.61	1.79
4	H	39	MET	SD-CE	-7.24	1.61	1.79
3	D	197	ILE	CG1-CD1	-6.86	1.25	1.51
3	E	145	MET	SD-CE	-6.16	1.64	1.79

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	38	A	C3'-C2'-O2'	7.62	126.04	114.60
5	I	38	A	C2'-C3'-O3'	6.63	119.44	109.50
3	E	65	VAL	CB-CA-C	5.80	115.75	110.13
3	C	185	ILE	N-CA-C	-5.78	106.15	112.80
5	I	21	U	C4'-C3'-O3'	-5.77	100.75	109.40

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	A	1897	0	1892	222	0
2	B	1709	0	1694	183	0
3	C	2488	0	2494	177	0
3	D	2503	0	2513	182	0
3	E	2450	0	2446	259	0
3	F	2450	0	2446	212	0
3	G	2131	0	2128	250	0
4	H	1800	0	1838	127	0
5	I	1317	0	659	191	0
6	J	722	0	421	39	0
7	K	759	0	418	77	0
8	L	4542	0	4629	193	0
9	A	1	0	0	0	0
All	All	24769	0	23578	1822	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 38.

The worst 5 of 1822 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:F:180:ARG:HG2	5:I:33:U:C5'	1.21	1.62
2:B:102:LEU:CD2	2:B:104:LEU:HD23	1.29	1.62
3:G:48:VAL:HG21	3:G:62:LEU:CD1	1.19	1.61
3:D:70:MET:CA	3:D:73:LEU:HD23	1.24	1.60
3:G:82:VAL:CG2	3:G:277:LEU:HD21	1.26	1.59

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	239/253 (94%)	226 (95%)	13 (5%)	0	100	100
2	B	217/220 (99%)	204 (94%)	12 (6%)	1 (0%)	24	56
3	C	325/348 (93%)	310 (95%)	14 (4%)	1 (0%)	36	65
3	D	327/348 (94%)	312 (95%)	15 (5%)	0	100	100
3	E	320/348 (92%)	310 (97%)	10 (3%)	0	100	100
3	F	320/348 (92%)	306 (96%)	14 (4%)	0	100	100
3	G	276/348 (79%)	263 (95%)	13 (5%)	0	100	100
4	H	232/268 (87%)	228 (98%)	4 (2%)	0	100	100
8	L	586/626 (94%)	572 (98%)	14 (2%)	0	100	100
All	All	2842/3107 (92%)	2731 (96%)	109 (4%)	2 (0%)	49	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	185	ASP
3	C	128	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	210/222 (95%)	205 (98%)	5 (2%)	43	61
2	B	176/177 (99%)	173 (98%)	3 (2%)	53	67

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	263/274 (96%)	261 (99%)	2 (1%)	73	76
3	D	264/274 (96%)	262 (99%)	2 (1%)	73	76
3	E	259/274 (94%)	254 (98%)	5 (2%)	50	65
3	F	259/274 (94%)	257 (99%)	2 (1%)	73	76
3	G	227/274 (83%)	219 (96%)	8 (4%)	32	54
4	H	191/215 (89%)	187 (98%)	4 (2%)	47	64
8	L	480/503 (95%)	480 (100%)	0	100	100
All	All	2329/2487 (94%)	2298 (99%)	31 (1%)	59	71

5 of 31 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	E	77	THR
4	H	26	LEU
3	F	282	MET
4	H	214	VAL
3	G	258	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 30 such sidechains are listed below:

Mol	Chain	Res	Type
3	D	187	ASN
8	L	472	ASN
3	E	68	ASN
8	L	583	GLN
8	L	112	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
5	I	60/61 (98%)	43 (71%)	14 (23%)

5 of 43 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	I	2	U
5	I	3	G

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Mol	Chain	Res	Type
5	I	4	A
5	I	6	C
5	I	7	G

5 of 14 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
5	I	34	G
5	I	38	A
5	I	48	C
5	I	46	C
5	I	47	C

5.4 Non-standard residues in protein, DNA, RNA chains

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
5	23G	I	61	5	24,29,30	1.20	2 (8%)	34,45,48	2.12	9 (26%)

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
5	23G	I	61	5	-	0/7/35/36	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	I	61	23G	C5-C4	3.20	1.47	1.38
5	I	61	23G	C6-N1	-2.46	1.34	1.38

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	61	23G	C5-C4-N3	-6.25	118.45	128.39
5	I	61	23G	C2-N3-C4	5.09	121.07	112.30
5	I	61	23G	N9-C4-N3	4.64	135.23	125.95
5	I	61	23G	OC1-PC-OC2	3.32	120.62	109.89
5	I	61	23G	C6-C5-N7	3.19	136.09	130.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	I	61	23G	9	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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.

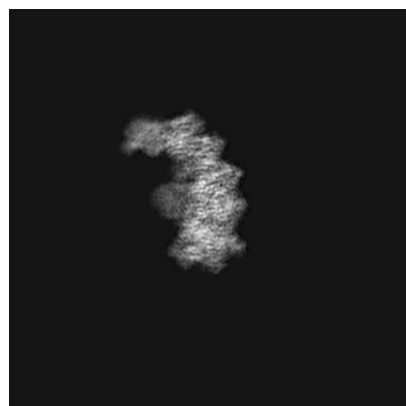
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-33185. These allow visual inspection of the internal detail of the map and identification of artifacts.

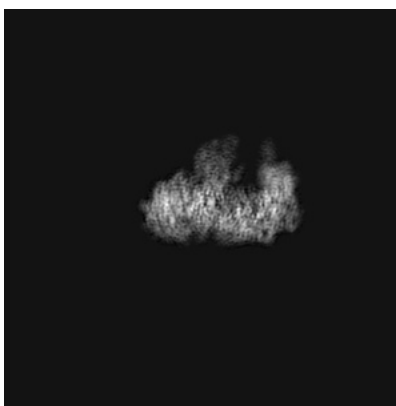
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

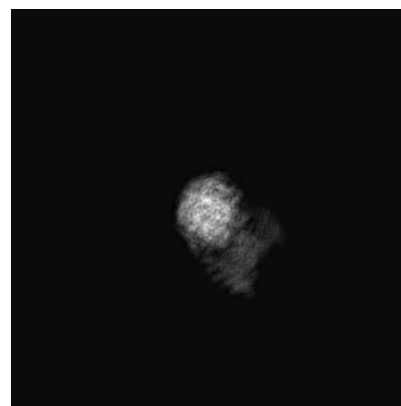
6.1.1 Primary map



X

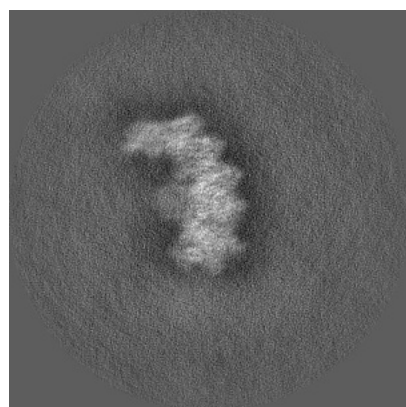


Y

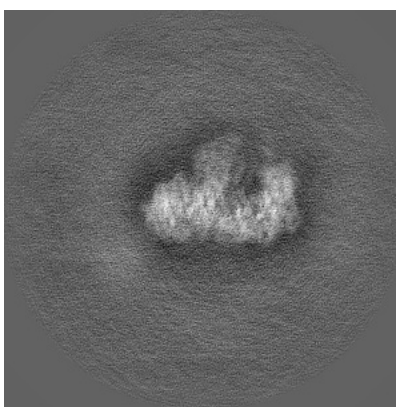


Z

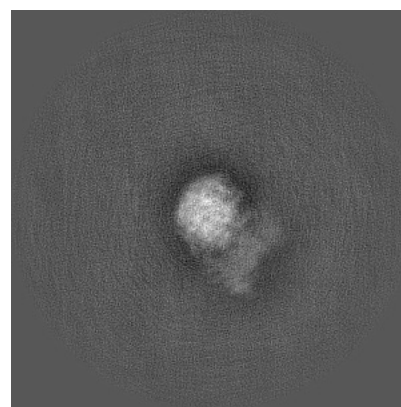
6.1.2 Raw map



X



Y

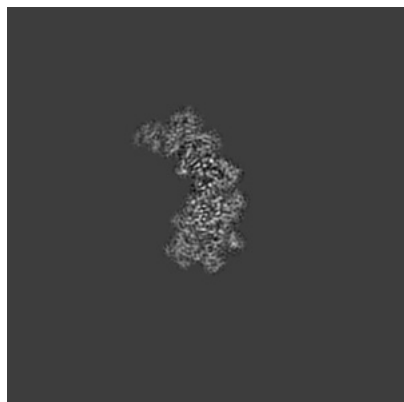


Z

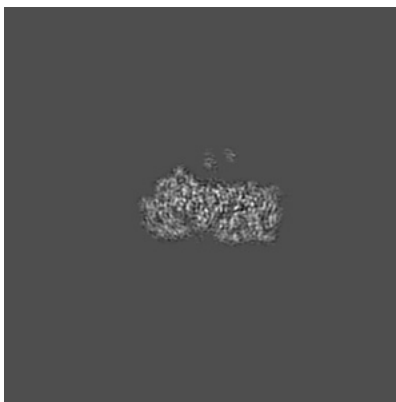
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 180

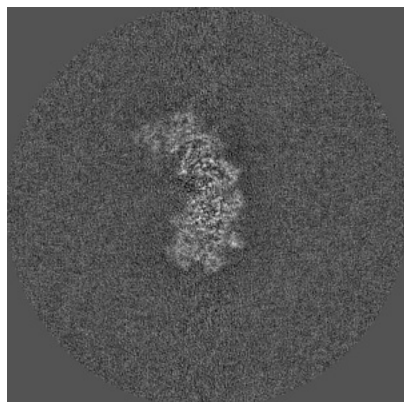


Y Index: 180

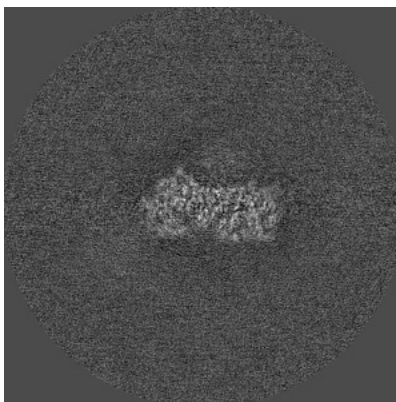


Z Index: 180

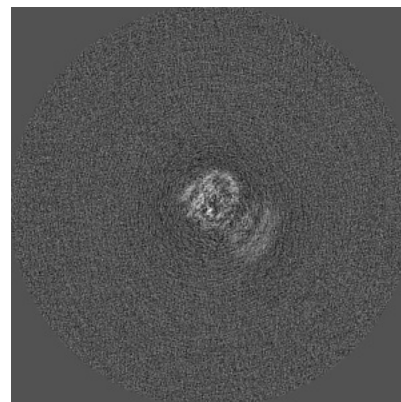
6.2.2 Raw map



X Index: 180



Y Index: 180

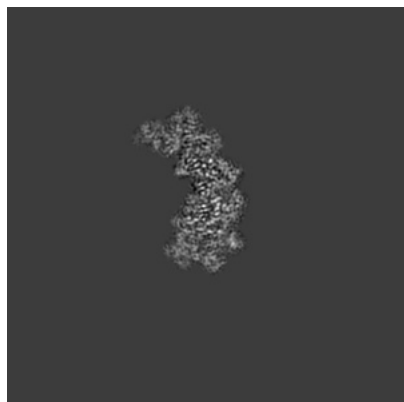


Z Index: 180

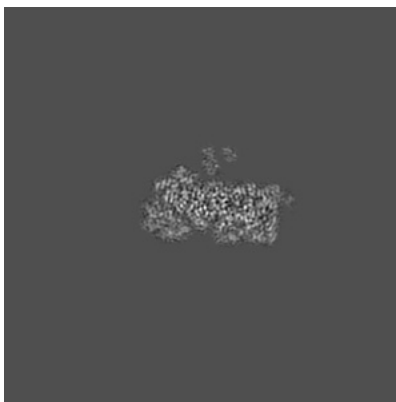
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 179

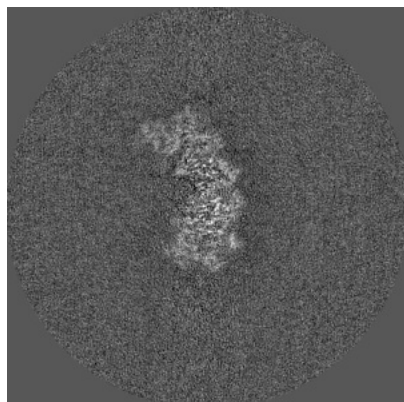


Y Index: 177

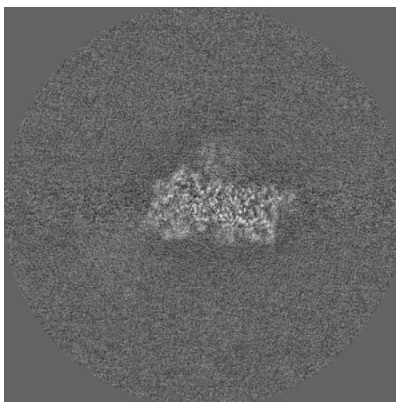


Z Index: 183

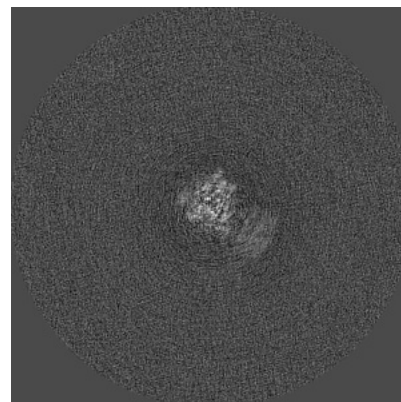
6.3.2 Raw map



X Index: 179



Y Index: 174

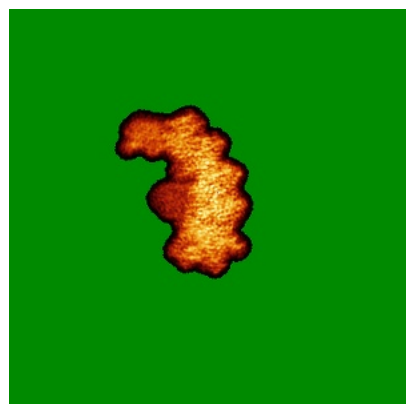


Z Index: 176

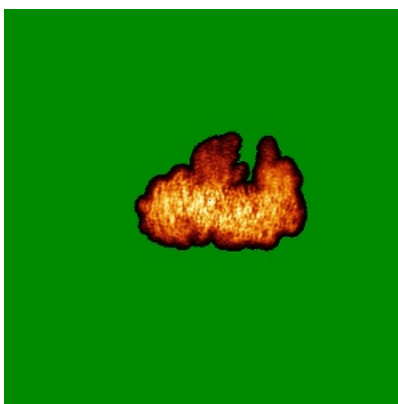
The images above show the largest variance slices of the map in three orthogonal directions.

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

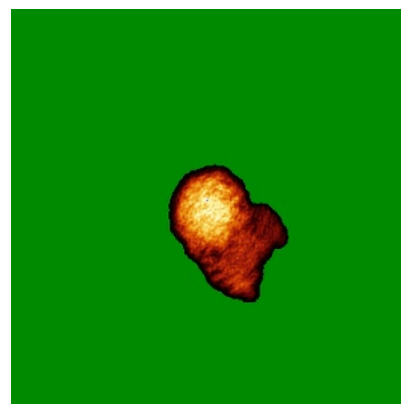
6.4.1 Primary map



X

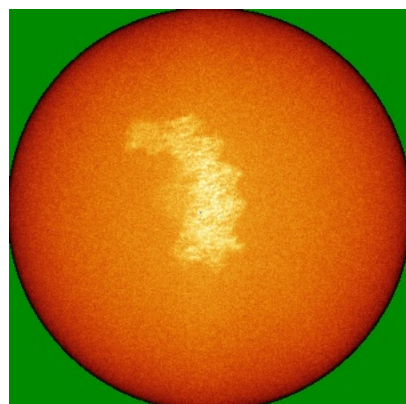


Y

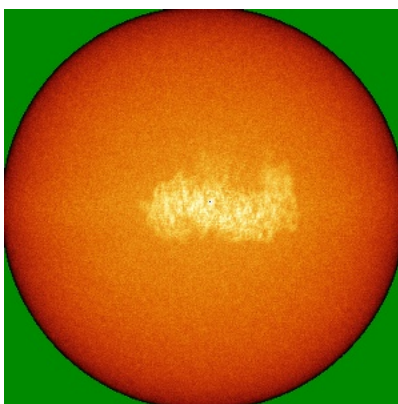


Z

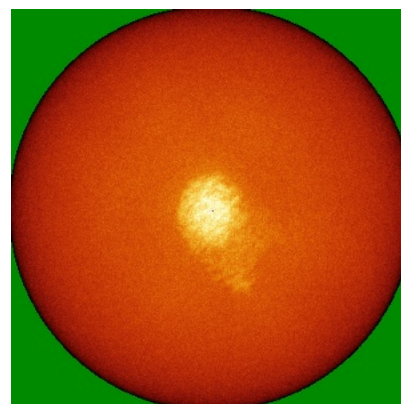
6.4.2 Raw map



X



Y



Z

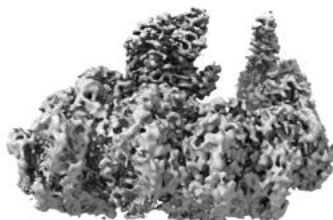
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

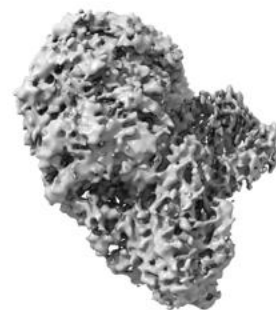
6.5.1 Primary map



X



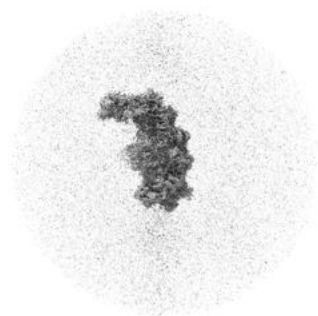
Y



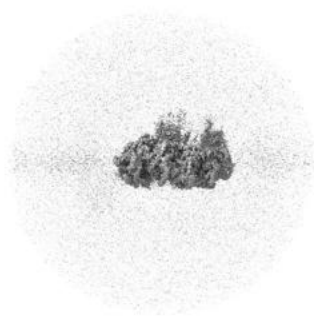
Z

The images above show the 3D surface view of the map at the recommended contour level 0.008. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X



Y



Z

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

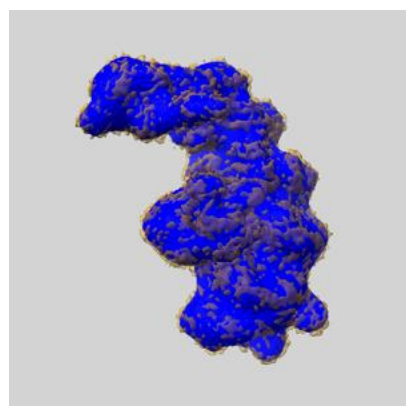
6.6 Mask visualisation [i](#)

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

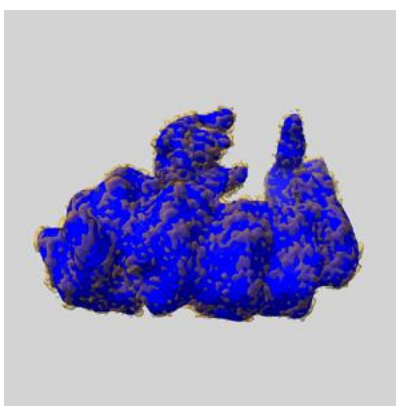
A mask typically either:

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

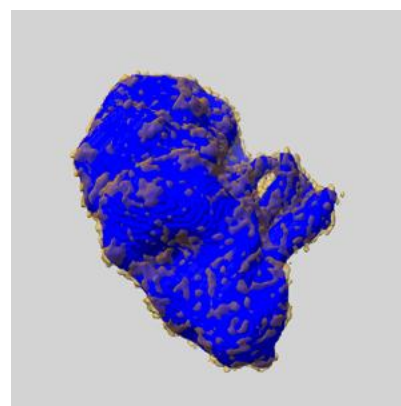
6.6.1 emd_33185_msk_1.map [i](#)



X



Y

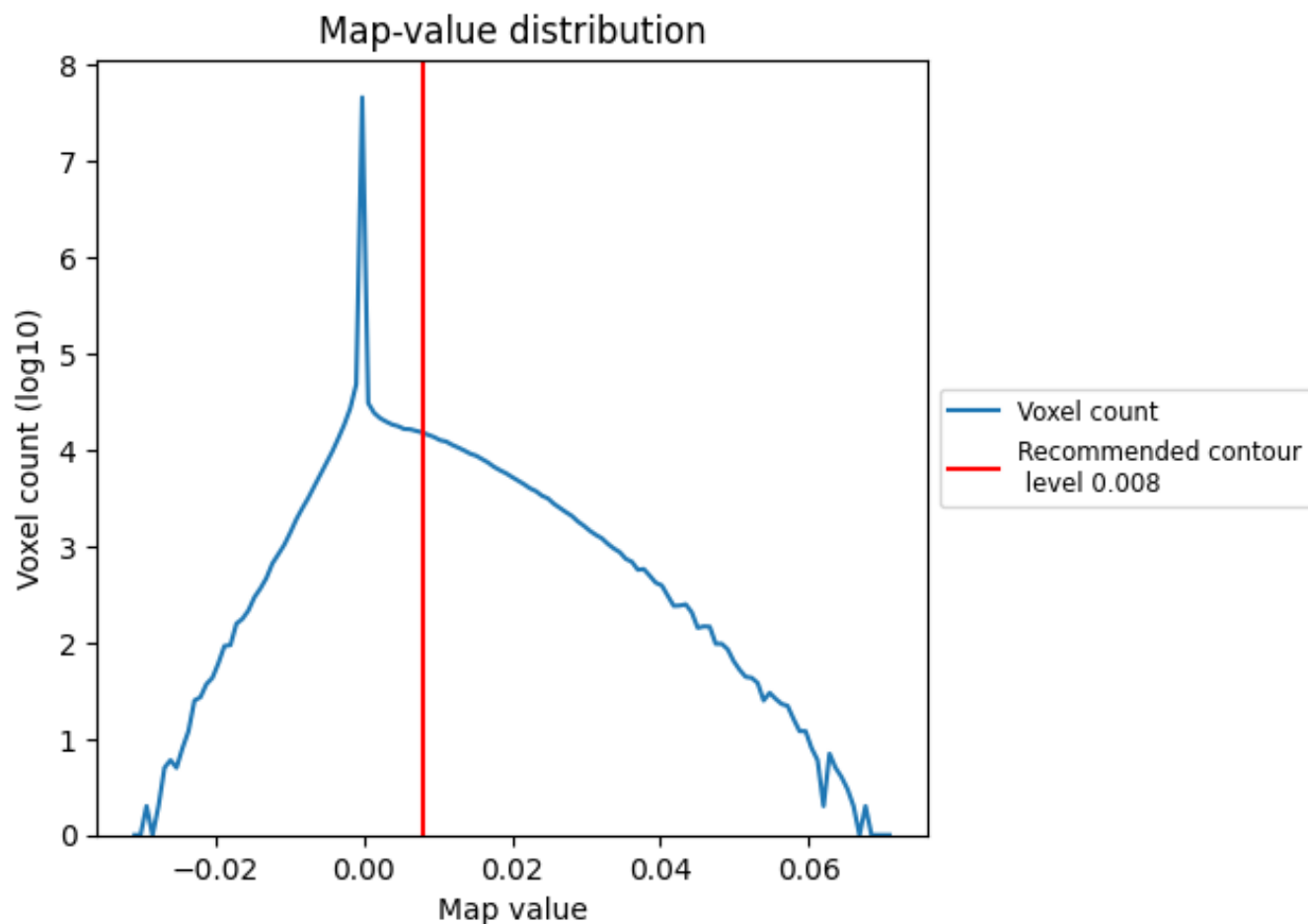


Z

7 Map analysis [i](#)

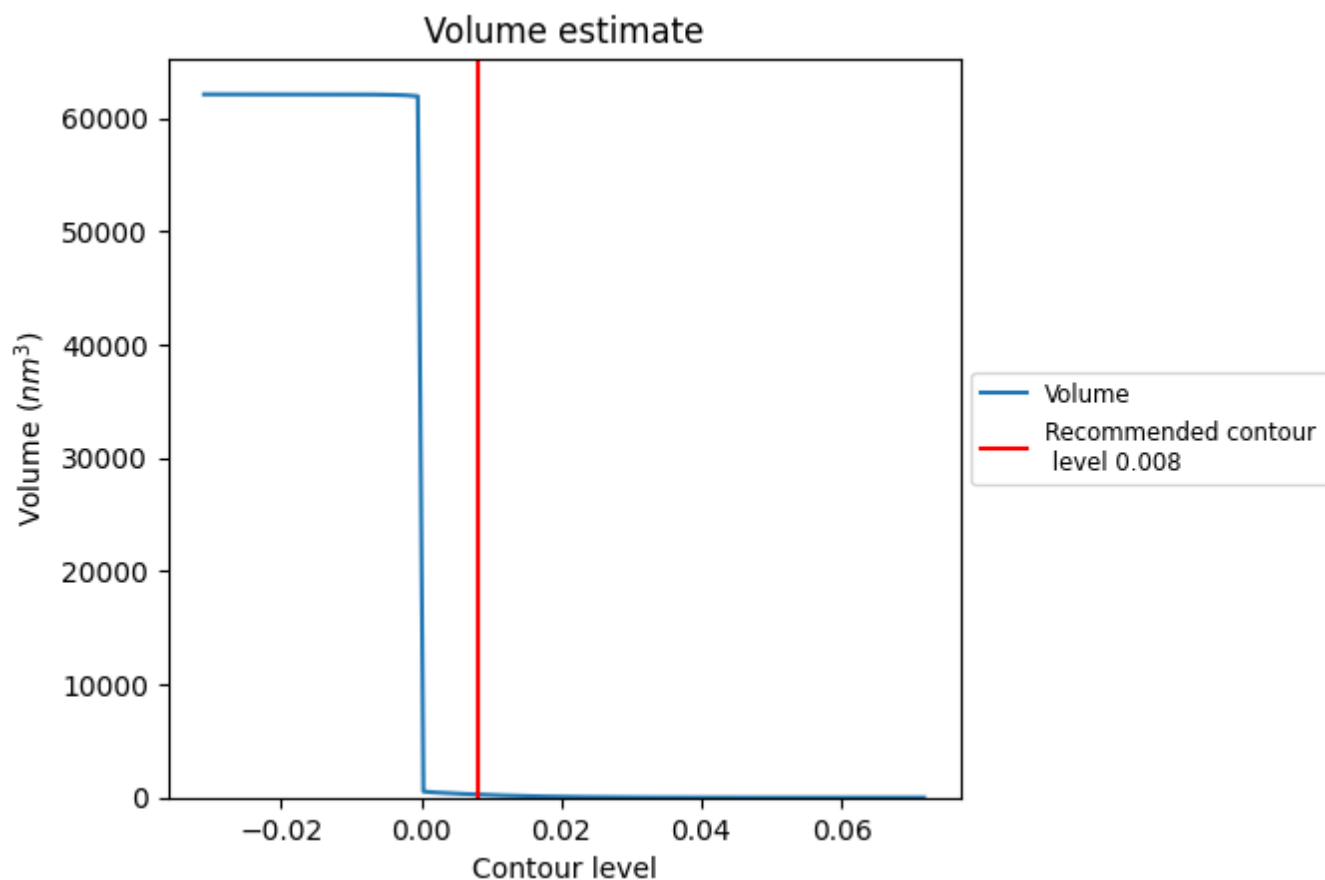
This section contains the results of statistical analysis of the map.

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.

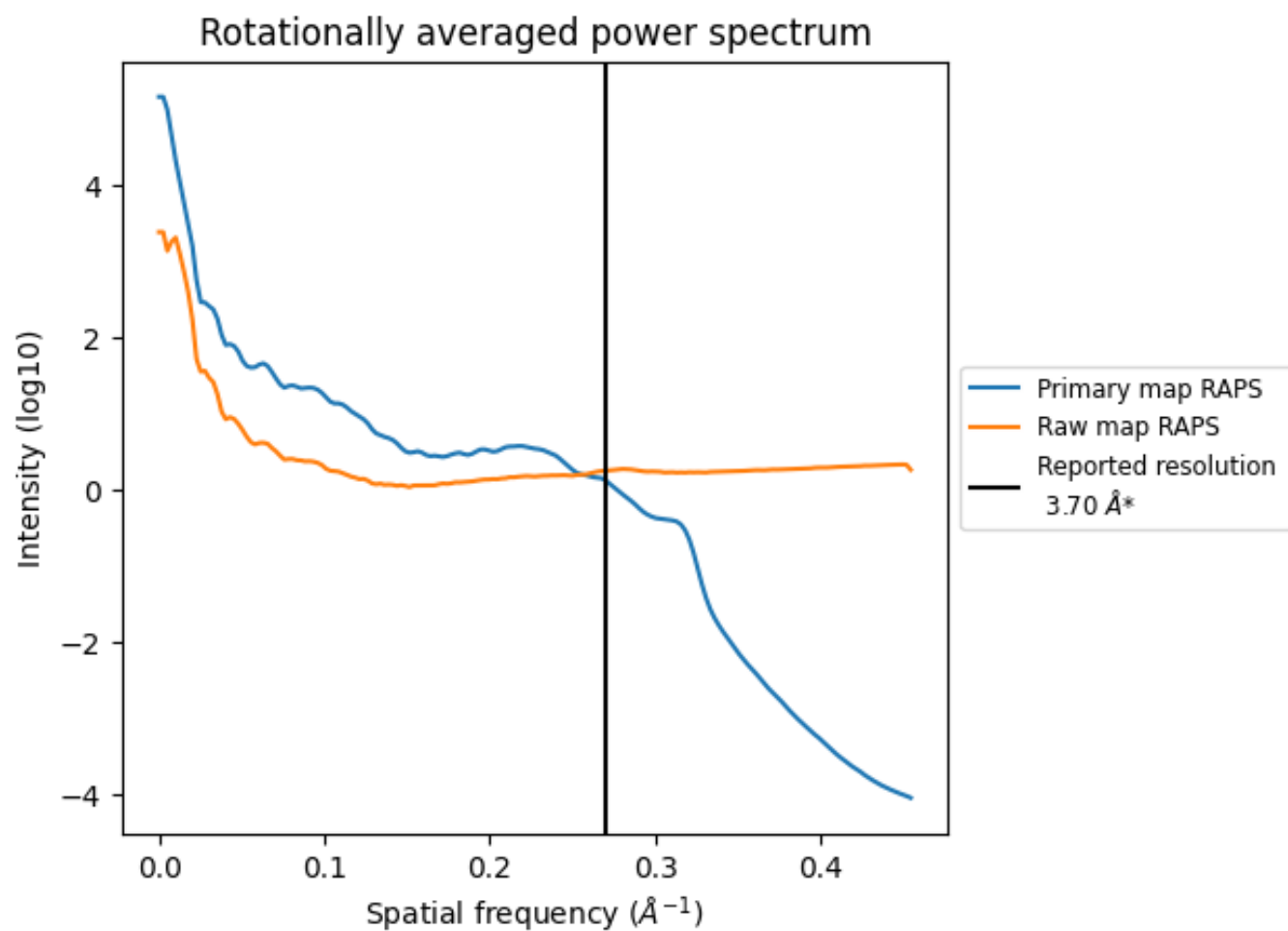
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 272 nm³; this corresponds to an approximate mass of 245 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 ⓘ

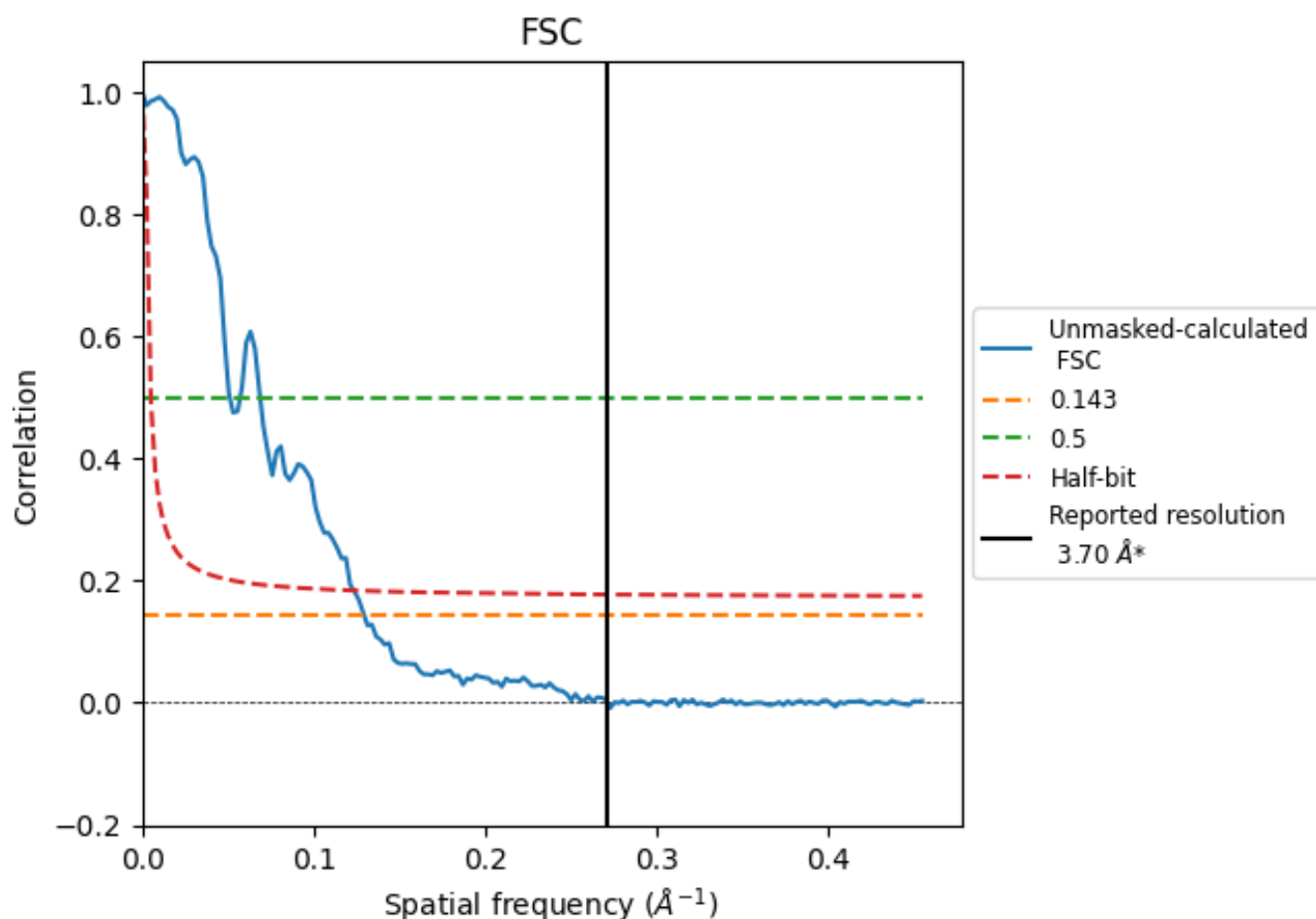


*Reported resolution corresponds to spatial frequency of 0.270 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.270 Å⁻¹

8.2 Resolution estimates [i](#)

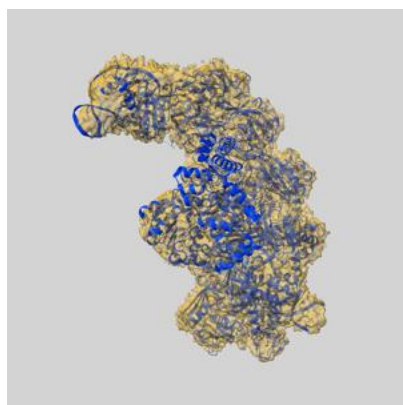
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	7.73	19.61	8.10

*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 7.73 differs from the reported value 3.7 by more than 10 %

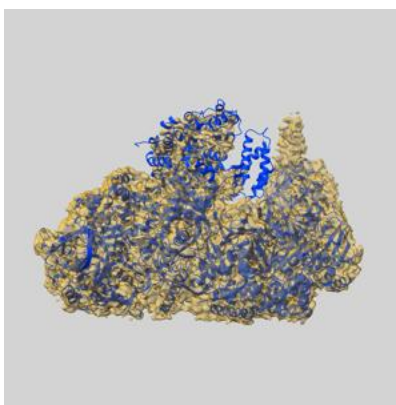
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-33185 and PDB model 7XG4. Per-residue inclusion information can be found in section [3](#) on page [6](#).

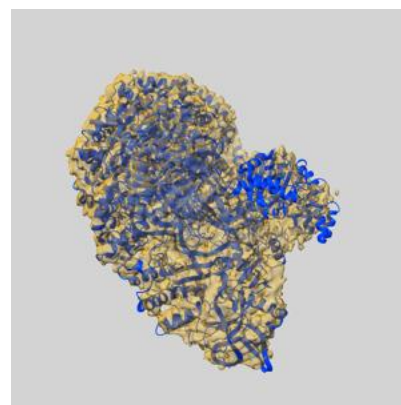
9.1 Map-model overlay [i](#)



X



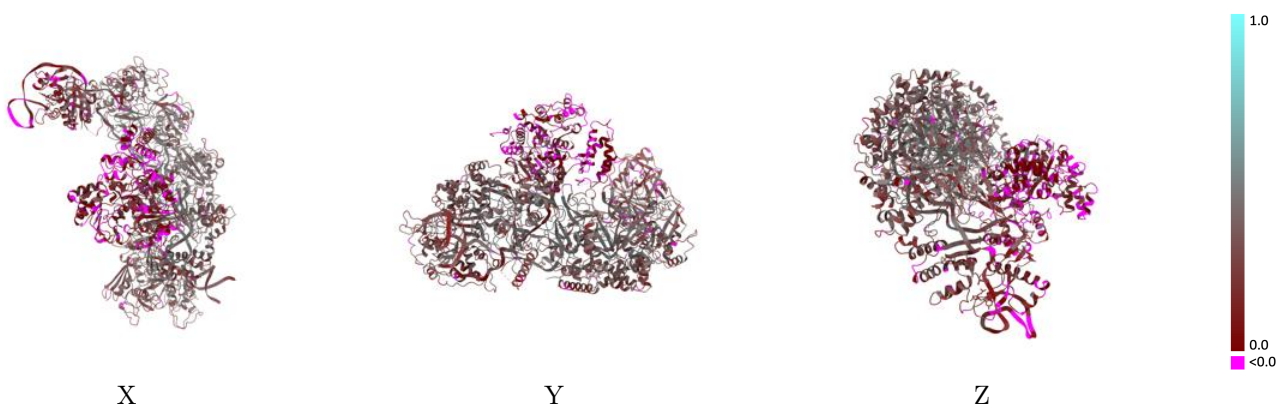
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.008 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](#)

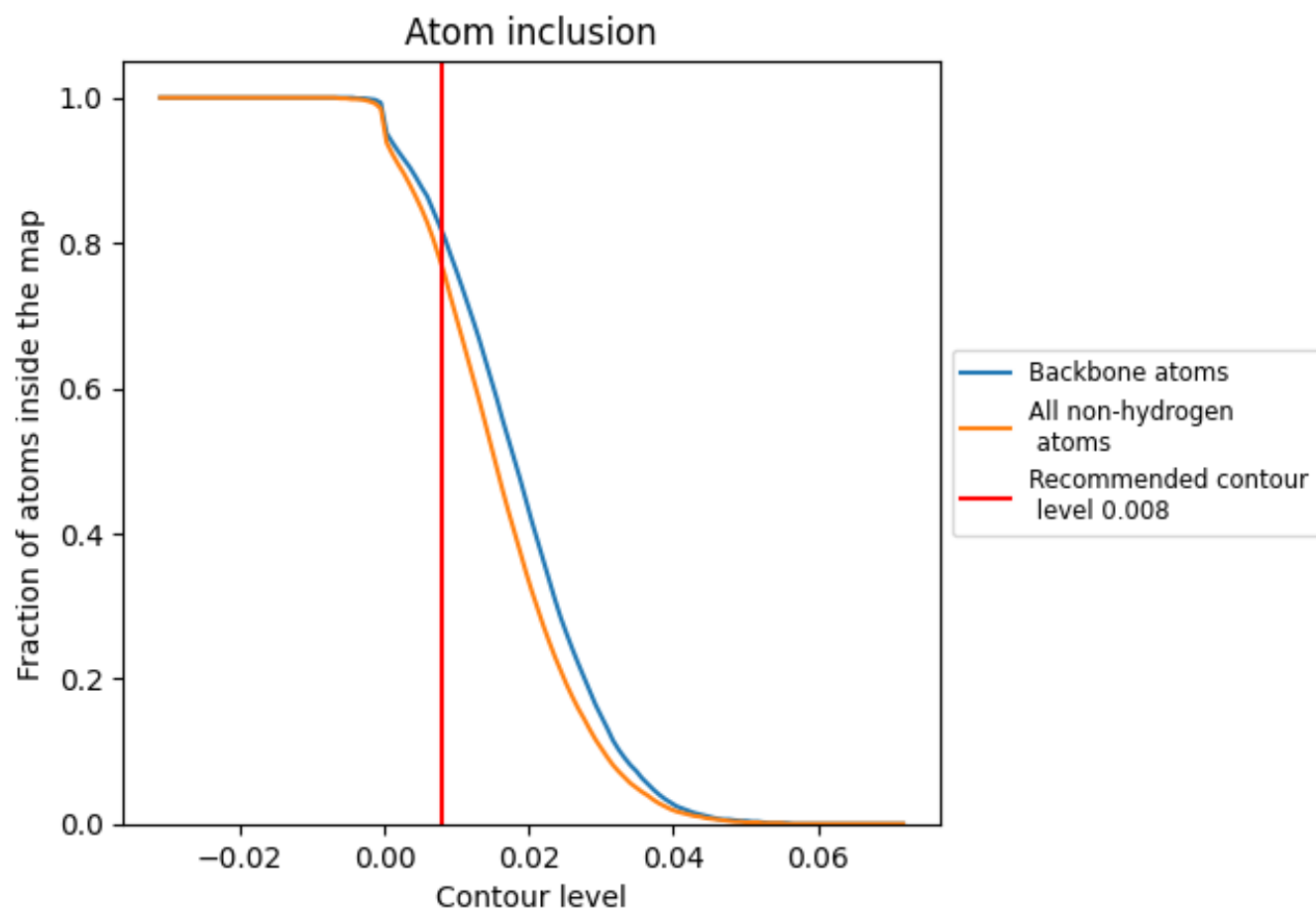


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 77% 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 (0.008) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7680	 0.2850
A	 0.8600	 0.3050
B	 0.8950	 0.3170
C	 0.9150	 0.3600
D	 0.9330	 0.3880
E	 0.9090	 0.3880
F	 0.8560	 0.3510
G	 0.8120	 0.3110
H	 0.7820	 0.2140
I	 0.8970	 0.3030
J	 0.6790	 0.1980
K	 0.9310	 0.3700
L	 0.3100	 0.0860

