



Full wwPDB EM Validation Report ⓘ

Apr 8, 2026 – 08:49 PM UTC

PDB ID : 9NZS / pdb_00009nzs
EMDB ID : EMD-49958
Title : Cryo-EM structure of OsCas12f-gRNA-DNA ternary complex. State III
Authors : Guan, K.; Taylor, D.W.
Deposited on : 2025-04-01
Resolution : 3.68 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:

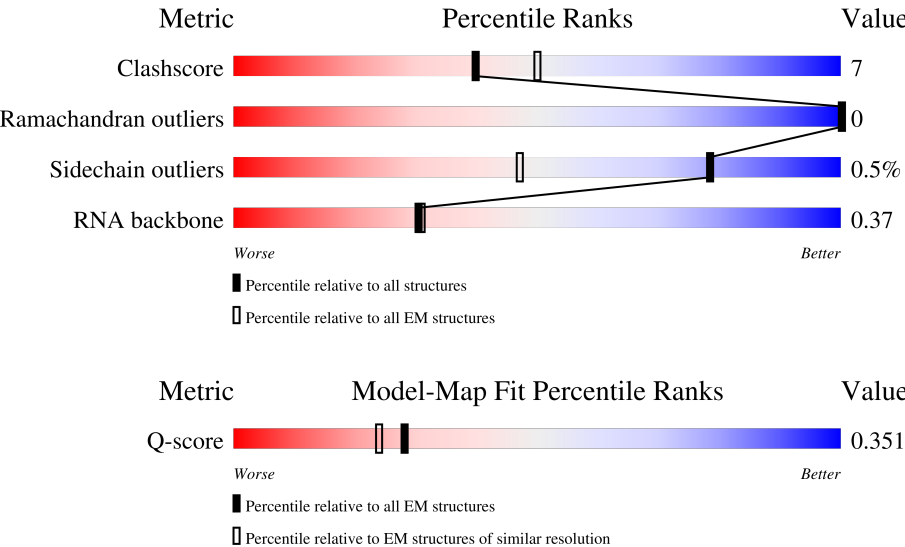
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.68 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	11376 (3.18 - 4.18)

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	433	
1	B	433	
2	C	45	

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Mol	Chain	Length	Quality of chain
3	D	45	<div><div></div><div>16%</div><div></div><div>82%</div></div>
4	E	151	<div><div></div><div>34%</div><div>44%</div><div>9%</div><div>13%</div></div>
5	G	4	<div><div></div><div>25%</div><div>75%</div><div>25%</div></div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9446 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 Cas12f.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	412	Total	C	N	O	S	0	0
			3295	2068	597	607	23		
1	B	316	Total	C	N	O	S	0	0
			2548	1617	449	467	15		

- Molecule 2 is a DNA chain called Targer strand DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	26	Total	C	N	O	P	0	0
			533	253	107	147	26		

- Molecule 3 is a DNA chain called Non-target strand DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	8	Total	C	N	O	P	0	0
			163	79	26	50	8		

- Molecule 4 is a RNA chain called guide RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	132	Total	C	N	O	P	0	0
			2830	1263	520	915	132		

- Molecule 5 is a DNA chain called DNA target poly C.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	G	4	Total	C	N	O	P	0	0
			76	36	12	24	4		

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

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

DA DG DC DA DA DG DG DC DA DA DC DG DC DT DC DC

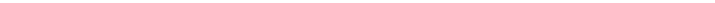
A17 G20 C23 C24 C25 A26 A27 G28 C42 DG DT DA

- Chain D: 16% 82%

DT	DA	G7		A10	C14	DT	DG	DC	DT	DT	DG	DT	DC	DT	DT	DG	DG	DG	DT	DG	DC	DA	DT	DT	DG	DG	DA	DG	DC	DT	DT	DG	DC	DC	DT	DT	DG	DC	DT
----	----	----	--	-----	-----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----	----

- Chain E: 

G	A75	A76	A77	A78	U81	A82	A83	A84	A85	U86	C90	A91	G96	U97	A98	A99	U100	C106	A107	U108	G109	A110	G111	C115	G116	A117	A118	A119	G120	C121	G122	G123	G127	A128	A129	G130	G133	C134	U137	C138	U139	A140	G141	G142	G143	U144	G145	C146	A147	U148	U149	G						
A	G2	A8	C19	A20	A21	A22	A23	U24	C25	G26	A27	G28	A29	C30	A31	G32	U33	A34	G35	C36	C37	G38	U39	A40	A41	A	A	C	G	U	U	U	G	A	G	U	U	U	C	A	A	G	C	G	U58	U59	G60	G61	G64	A65	C66	A67	C68	A69	C70	U71	G72	G73

- Chain G: 

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	42153	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS GLACIOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	49	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	0.466	Depositor
Minimum map value	-0.186	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.013	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	300.8, 300.8, 300.8	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.93999994, 0.93999994, 0.93999994	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.33	0/3359	0.73	1/4514 (0.0%)
1	B	0.31	0/2596	0.76	4/3488 (0.1%)
2	C	0.32	0/600	0.57	0/922
3	D	0.29	0/181	0.57	0/277
4	E	0.24	0/3167	0.55	2/4935 (0.0%)
5	G	0.33	0/83	0.59	0/124
All	All	0.30	0/9986	0.67	7/14260 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	20	A	OP1-P-O3'	-9.65	79.03	108.00
1	B	214	PRO	CA-N-CD	-7.27	101.82	112.00
1	B	120	MET	CB-CG-SD	6.39	131.87	112.70
1	B	120	MET	CA-CB-CG	6.37	126.84	114.10
4	E	20	A	OP2-P-O3'	-6.24	89.27	108.00
1	A	213	PRO	CA-N-CD	-5.55	104.23	112.00
1	B	120	MET	CB-CA-C	5.43	118.49	109.53

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3295	0	3281	45	0
1	B	2548	0	2547	43	0
2	C	533	0	290	6	0
3	D	163	0	93	1	0
4	E	2830	0	1428	38	0
5	G	76	0	45	1	0
6	A	1	0	0	0	0
All	All	9446	0	7684	121	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 7.

All (121) 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:E:34:A:H2	4:E:64:G:H1	1.23	0.83
1:A:267:MET:HB3	1:A:290:VAL:HG22	1.64	0.79
1:A:307:ARG:HE	4:E:129:A:H8	1.32	0.78
1:A:31:TRP:HE1	1:A:200:LYS:HG2	1.51	0.73
2:C:17:DA:H61	4:E:149:U:H3	1.36	0.72
1:B:120:MET:HE1	1:B:123:PRO:HG3	1.80	0.63
1:B:137:THR:HG22	1:B:153:LEU:HD23	1.81	0.62
1:B:219:LEU:HG	1:B:221:PRO:HD3	1.82	0.61
1:B:17:LEU:HD12	1:B:18:ASP:H	1.66	0.61
1:B:229:MET:HG3	1:B:235:LEU:HD13	1.82	0.60
1:B:33:LEU:HD12	1:B:131:LEU:HD11	1.83	0.59
1:A:179:ARG:HH12	4:E:8:A:H1'	1.66	0.59
1:B:137:THR:HB	1:B:153:LEU:HB3	1.86	0.58
1:B:140:LEU:HD22	1:B:150:THR:HG22	1.86	0.57
1:B:196:GLN:HE21	1:B:208:VAL:HG12	1.69	0.57
1:B:235:LEU:HB3	1:B:248:ILE:HB	1.87	0.56
1:A:42:ASN:HD21	1:A:125:TYR:HB2	1.70	0.56
1:A:145:ARG:HB3	1:A:147:PRO:HD2	1.88	0.56
4:E:139:U:H2'	4:E:140:U:C6	2.42	0.55
1:B:138:VAL:HG22	1:B:152:THR:HA	1.88	0.54
1:A:157:LYS:HD2	3:D:10:DA:H4'	1.89	0.54
1:A:159:LYS:HZ2	1:A:166:ASN:HD21	1.56	0.53
1:B:31:TRP:HH2	1:B:200:LYS:HE3	1.73	0.53
4:E:140:U:H2'	4:E:141:G:H8	1.74	0.53
1:A:415:ILE:HG23	1:A:418:ILE:HB	1.92	0.52
4:E:34:A:H2	4:E:64:G:N1	2.01	0.52
1:B:312:LEU:HD12	1:B:313:ILE:HG12	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:390:SER:HB3	1:A:393:ASP:HB2	1.92	0.51
1:B:235:LEU:HD21	1:B:312:LEU:HD23	1.92	0.50
4:E:26:G:H2'	4:E:27:A:H8	1.75	0.50
4:E:33:U:H2'	4:E:34:A:C8	2.46	0.50
1:A:291:TYR:CE1	2:C:25:DC:H4'	2.46	0.50
4:E:34:A:N1	4:E:64:G:O6	2.45	0.50
4:E:138:C:H2'	4:E:139:U:C6	2.47	0.50
1:B:160:ARG:HH12	1:B:168:LYS:HZ3	1.59	0.50
1:A:9:VAL:HG12	1:A:209:THR:HG22	1.94	0.50
1:A:38:ARG:HH12	1:A:127:MET:HA	1.77	0.50
1:A:377:CYS:HB2	1:A:396:CYS:SG	2.51	0.49
4:E:109:G:H2'	4:E:110:A:H8	1.77	0.49
4:E:140:U:H2'	4:E:141:G:C8	2.48	0.49
4:E:110:A:H2'	4:E:111:G:H8	1.77	0.49
1:A:181:ILE:O	1:A:185:LEU:HB2	2.13	0.49
1:A:15:ARG:HH21	4:E:100:U:H1'	1.78	0.49
1:B:92:MET:HE2	1:B:92:MET:HB3	1.70	0.48
4:E:110:A:H2'	4:E:111:G:C8	2.48	0.48
1:B:418:ILE:O	1:B:422:ILE:HG22	2.14	0.47
1:B:227:VAL:HG23	1:B:237:ALA:HB2	1.95	0.47
4:E:109:G:H2'	4:E:110:A:C8	2.49	0.47
4:E:143:G:O2'	4:E:144:U:H5'	2.14	0.47
1:A:142:GLU:HA	1:A:148:ILE:HG22	1.96	0.47
1:A:272:ALA:HB2	4:E:144:U:H5''	1.96	0.47
1:B:28:GLU:HA	1:B:31:TRP:CD1	2.49	0.47
1:A:17:LEU:HD11	1:A:26:MET:HB3	1.96	0.47
1:B:304:ILE:HD11	1:B:308:TYR:CZ	2.50	0.47
1:A:22:ASP:HB2	1:A:25:ASN:HB3	1.97	0.46
4:E:26:G:H2'	4:E:27:A:C8	2.50	0.46
4:E:141:G:H2'	4:E:142:G:H8	1.79	0.46
1:B:185:LEU:HD12	1:B:190:TYR:HB2	1.97	0.46
4:E:133:G:H2'	4:E:134:C:C6	2.50	0.46
1:A:290:VAL:HG12	1:A:294:LYS:HE2	1.96	0.46
1:A:269:LYS:HE2	1:A:269:LYS:HB2	1.78	0.46
1:A:38:ARG:NH1	1:A:127:MET:HA	2.30	0.46
1:A:130:PRO:HB3	1:A:199:TYR:CD1	2.50	0.46
1:B:15:ARG:HB2	1:B:205:PHE:HB3	1.96	0.46
1:B:216:GLU:HG2	1:B:218:PRO:HD3	1.98	0.46
1:A:377:CYS:HA	1:A:408:ASN:HD21	1.81	0.46
1:A:17:LEU:HD21	1:A:26:MET:HG3	1.97	0.46
1:A:235:LEU:HD13	1:A:350:LEU:HD21	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:144:U:H2'	4:E:145:G:C8	2.51	0.45
1:B:255:LYS:HD2	4:E:22:A:C5	2.51	0.45
4:E:137:U:H2'	4:E:138:C:C6	2.51	0.45
1:B:168:LYS:HB3	1:B:168:LYS:HE3	1.74	0.45
1:B:312:LEU:HD11	1:B:354:ILE:HG12	1.99	0.45
1:A:130:PRO:HB3	1:A:199:TYR:HD1	1.81	0.44
1:B:15:ARG:HG2	2:C:20:DG:H5''	2.00	0.44
1:B:360:GLU:HG3	1:B:361:HIS:CD2	2.52	0.44
4:E:90:C:H2'	4:E:91:A:H8	1.83	0.44
1:A:114:GLU:HB3	1:A:120:MET:HE3	1.98	0.44
1:B:224:ILE:HD13	1:B:415:ILE:HG22	2.00	0.44
4:E:106:C:H2'	4:E:107:A:C8	2.52	0.44
4:E:115:C:H2'	4:E:116:G:C4	2.53	0.44
4:E:106:C:H2'	4:E:107:A:H8	1.82	0.44
1:B:201:ARG:HB2	1:B:203:LYS:O	2.18	0.43
1:B:96:ASN:O	1:B:100:THR:HG22	2.18	0.43
1:A:27:GLN:HG2	1:A:31:TRP:CH2	2.54	0.43
4:E:96:G:H4'	4:E:97:U:C2	2.53	0.43
1:A:14:LEU:HD23	1:A:171:MET:SD	2.59	0.43
1:B:17:LEU:HD22	1:B:206:LEU:HB2	2.01	0.43
1:A:384:ASP:HB2	1:A:387:ASN:HB2	2.00	0.43
1:B:227:VAL:HG22	1:B:235:LEU:HD11	1.99	0.43
1:B:41:LEU:HD23	1:B:104:ALA:HB2	2.02	0.42
1:A:49:PHE:HE2	1:B:122:ILE:HG21	1.84	0.42
1:A:185:LEU:HD22	1:A:190:TYR:HB3	2.00	0.42
1:A:404:ASN:HD22	1:A:407:PHE:HB2	1.83	0.42
1:B:16:TYR:HB2	1:B:174:HIS:HB3	2.00	0.42
2:C:26:DA:H2''	2:C:27:DA:H5'	2.01	0.42
4:E:73:G:H2'	4:E:74:A:H8	1.84	0.42
4:E:149:U:H5''	4:E:149:U:H6	1.84	0.42
1:B:116:LEU:HD23	1:B:116:LEU:HA	1.82	0.42
1:B:297:VAL:O	1:B:301:ARG:HG2	2.19	0.42
1:A:416:ARG:HD3	1:A:417:ASN:N	2.35	0.42
1:B:225:LEU:HB2	1:B:321:CYS:SG	2.60	0.42
1:A:194:GLU:HB3	1:A:209:THR:OG1	2.20	0.42
1:B:16:TYR:HD2	1:B:179:ARG:HH11	1.68	0.42
2:C:23:DC:H2'	2:C:24:DC:C6	2.55	0.42
1:A:255:LYS:HB2	1:A:255:LYS:HE2	1.90	0.41
1:B:160:ARG:HH12	1:B:168:LYS:NZ	2.18	0.41
4:E:36:C:C4	4:E:37:C:N4	2.88	0.41
1:A:290:VAL:O	1:A:294:LYS:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:ALA:H	5:G:4:DC:H5''	1.86	0.41
1:A:277:GLY:HA3	4:E:24:U:H5''	2.03	0.41
4:E:2:G:H22	4:E:19:C:N4	2.19	0.41
1:A:31:TRP:NE1	1:A:200:LYS:HG2	2.27	0.40
1:B:31:TRP:HZ3	1:B:200:LYS:HG2	1.85	0.40
4:E:34:A:N1	4:E:64:G:C6	2.89	0.40
1:A:239:THR:HG23	1:A:242:GLU:HB2	2.03	0.40
4:E:122:G:H2'	4:E:123:G:C8	2.57	0.40
1:A:225:LEU:HD23	1:A:324:ILE:HG12	2.03	0.40
4:E:90:C:H2'	4:E:91:A:C8	2.56	0.40
1:A:346:THR:HG22	2:C:28:DG:H3'	2.03	0.40
1:B:418:ILE:HA	1:B:421:ILE:HG12	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	408/433 (94%)	395 (97%)	13 (3%)	0	100	100
1	B	304/433 (70%)	300 (99%)	4 (1%)	0	100	100
All	All	712/866 (82%)	695 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	353/370 (95%)	351 (99%)	2 (1%)	78	79
1	B	274/370 (74%)	273 (100%)	1 (0%)	84	81
All	All	627/740 (85%)	624 (100%)	3 (0%)	78	80

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	326	MET
1	A	327	GLU
1	B	364	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	50	GLN
1	A	166	ASN
1	A	268	GLN
1	A	273	HIS
1	A	344	HIS
1	A	361	HIS
1	A	404	ASN
1	A	412	ASN
1	B	29	GLN
1	B	34	GLN
1	B	50	GLN
1	B	102	GLN
1	B	128	ASN
1	B	351	GLN
1	B	412	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	E	131/151 (86%)	48 (36%)	8 (6%)

All (48) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	E	8	A
4	E	19	C
4	E	20	A
4	E	21	A
4	E	23	A
4	E	28	G
4	E	29	A
4	E	30	C
4	E	32	G
4	E	35	G
4	E	37	C
4	E	39	U
4	E	40	A
4	E	41	A
4	E	59	U
4	E	61	G
4	E	65	A
4	E	66	C
4	E	67	A
4	E	68	C
4	E	69	A
4	E	70	C
4	E	71	U
4	E	76	A
4	E	78	G
4	E	81	U
4	E	82	A
4	E	83	A
4	E	85	A
4	E	86	U
4	E	96	G
4	E	97	U
4	E	99	A
4	E	115	C
4	E	116	G
4	E	117	A
4	E	118	A
4	E	119	A
4	E	120	G
4	E	128	A
4	E	129	A
4	E	130	G
4	E	140	U

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
4	E	144	U
4	E	145	G
4	E	146	C
4	E	148	U
4	E	149	U

All (8) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	E	28	G
4	E	29	A
4	E	58	G
4	E	67	A
4	E	98	A
4	E	118	A
4	E	127	G
4	E	139	U

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](#)

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.

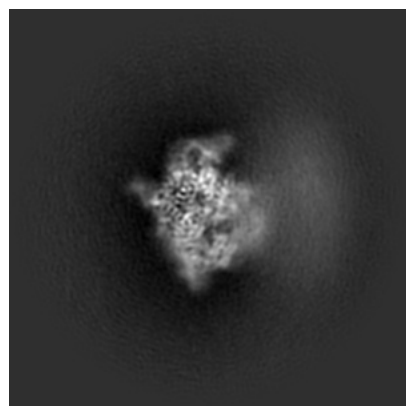
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-49958. These allow visual inspection of the internal detail of the map and identification of artifacts.

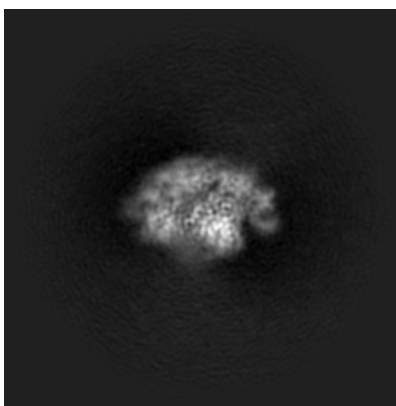
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

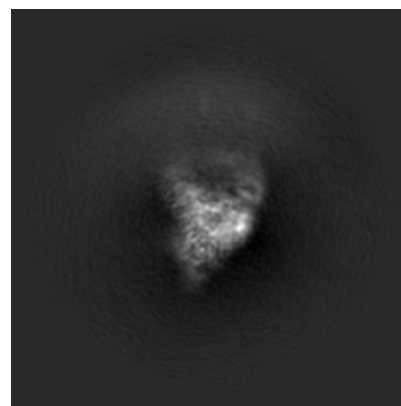
6.1.1 Primary map



X

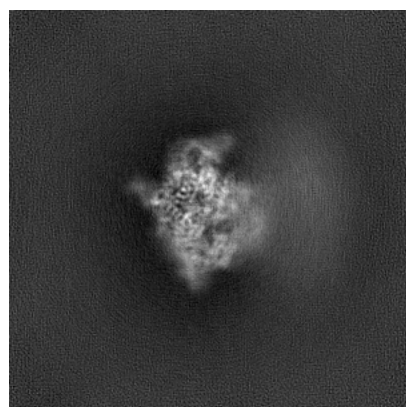


Y

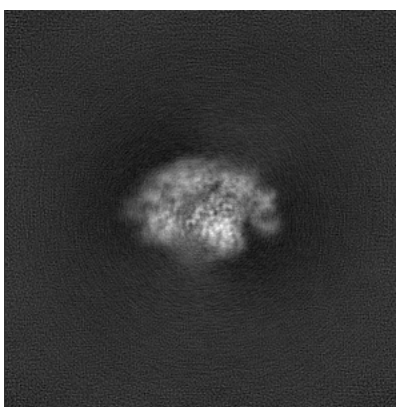


Z

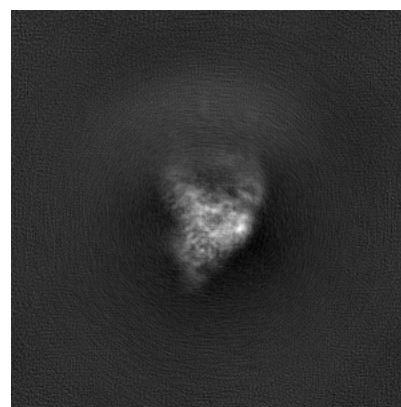
6.1.2 Raw map



X



Y

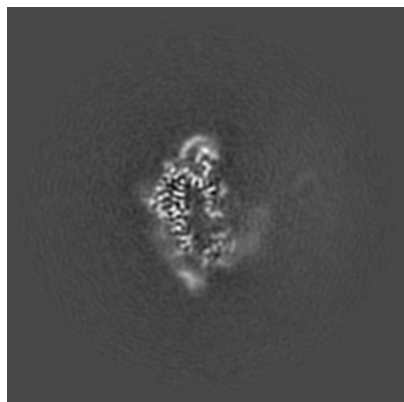


Z

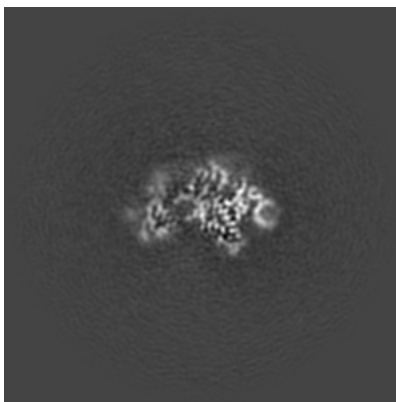
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

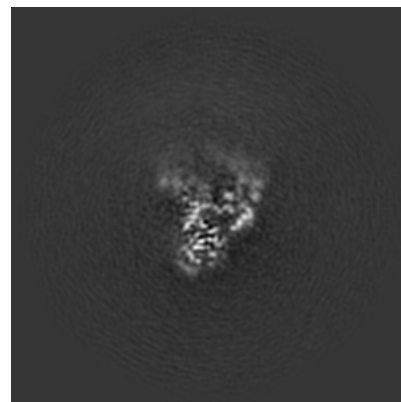
6.2.1 Primary map



X Index: 160

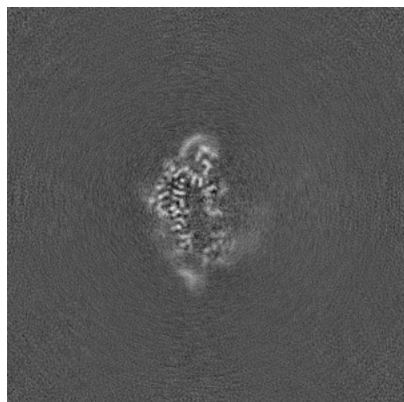


Y Index: 160

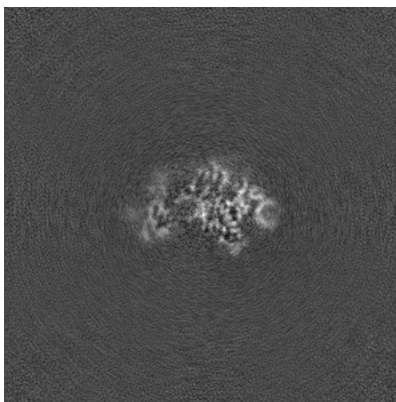


Z Index: 160

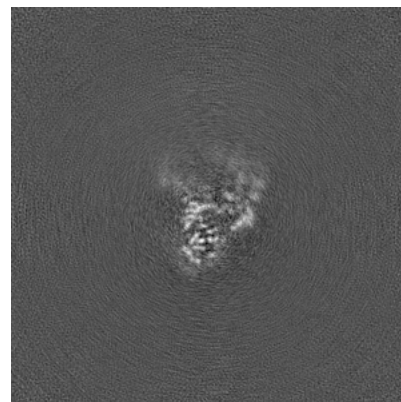
6.2.2 Raw map



X Index: 160



Y Index: 160

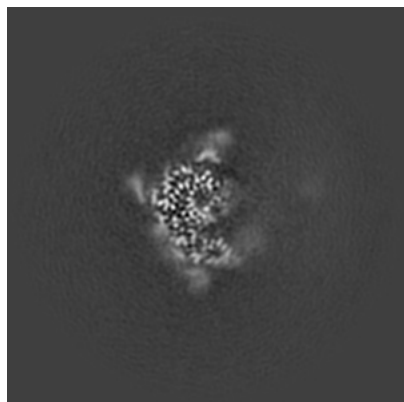


Z Index: 160

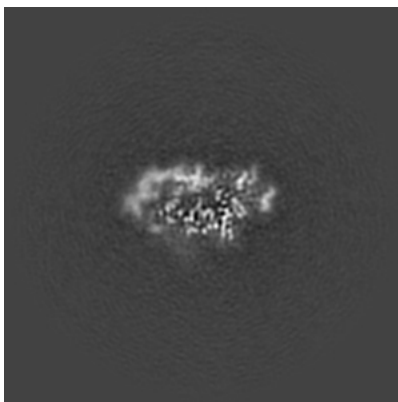
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

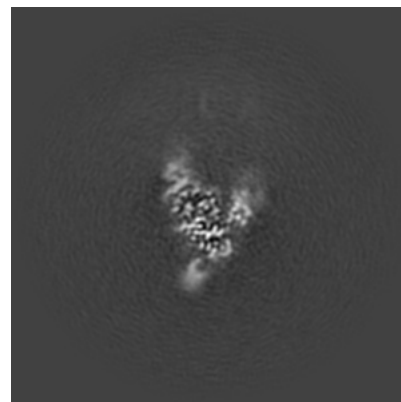
6.3.1 Primary map



X Index: 152

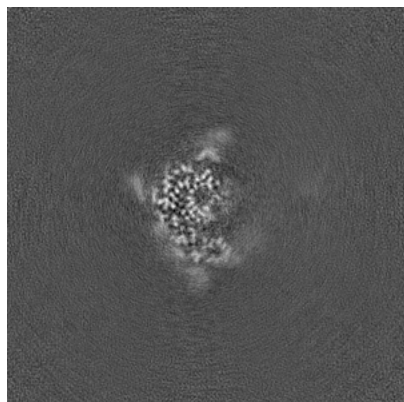


Y Index: 144

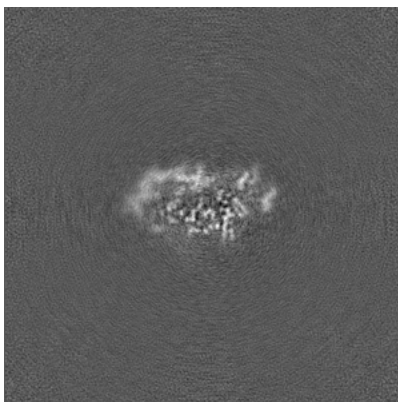


Z Index: 175

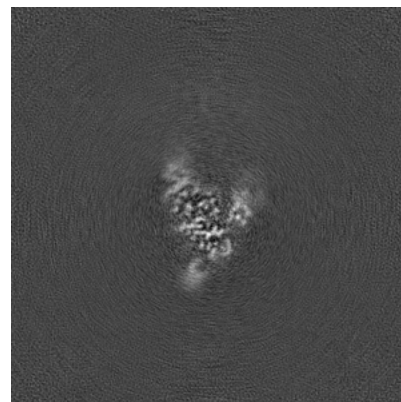
6.3.2 Raw map



X Index: 152



Y Index: 144

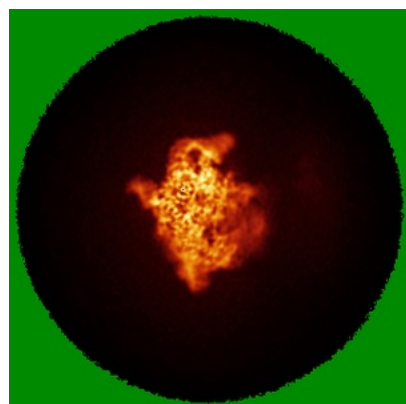


Z Index: 176

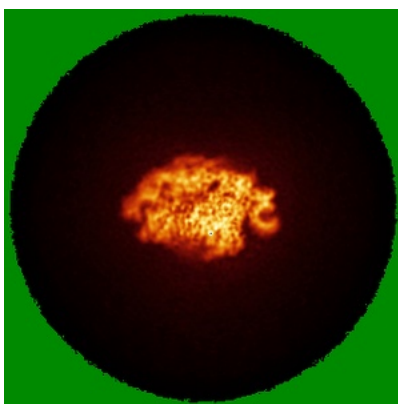
The images above show the largest variance slices of the map in three orthogonal directions.

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

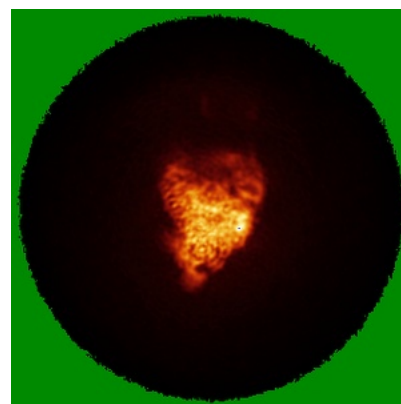
6.4.1 Primary map



X

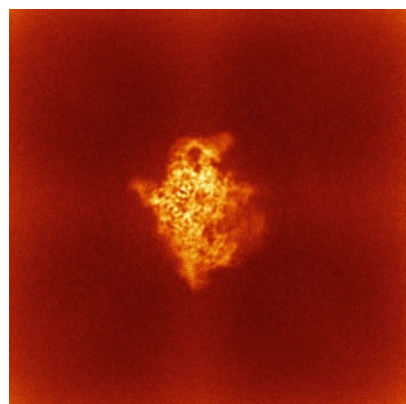


Y

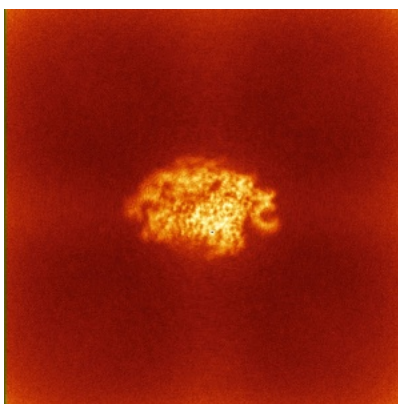


Z

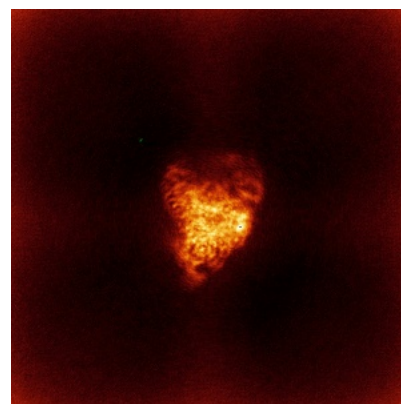
6.4.2 Raw map



X



Y

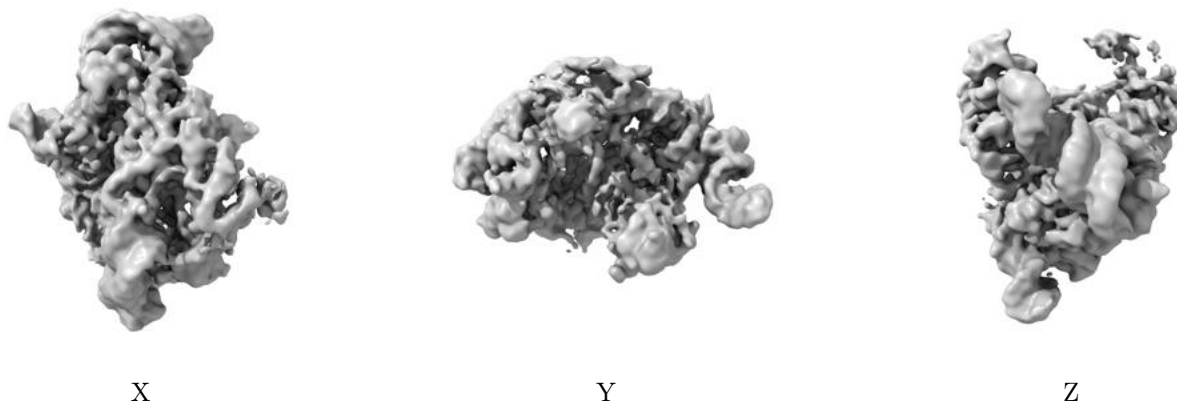


Z

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

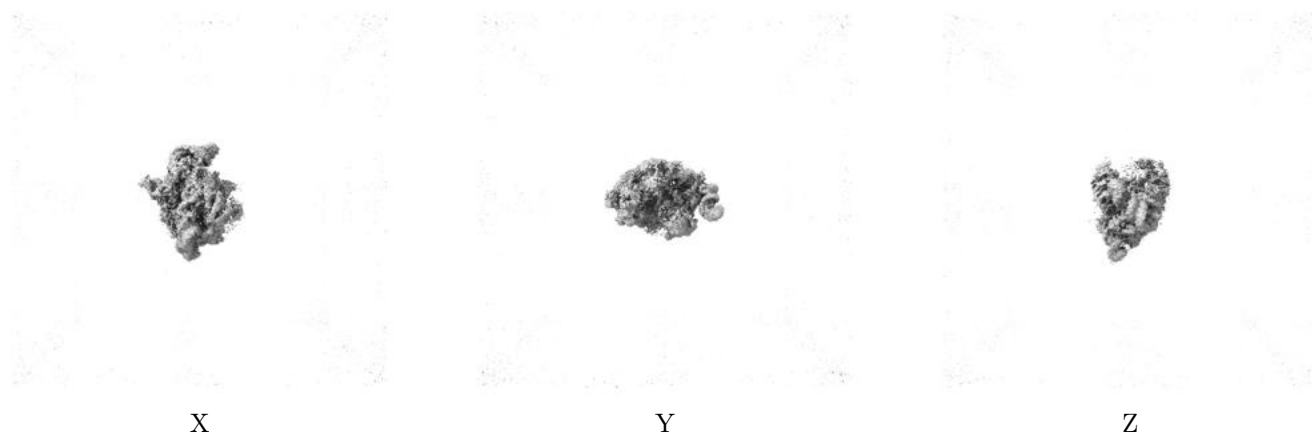
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

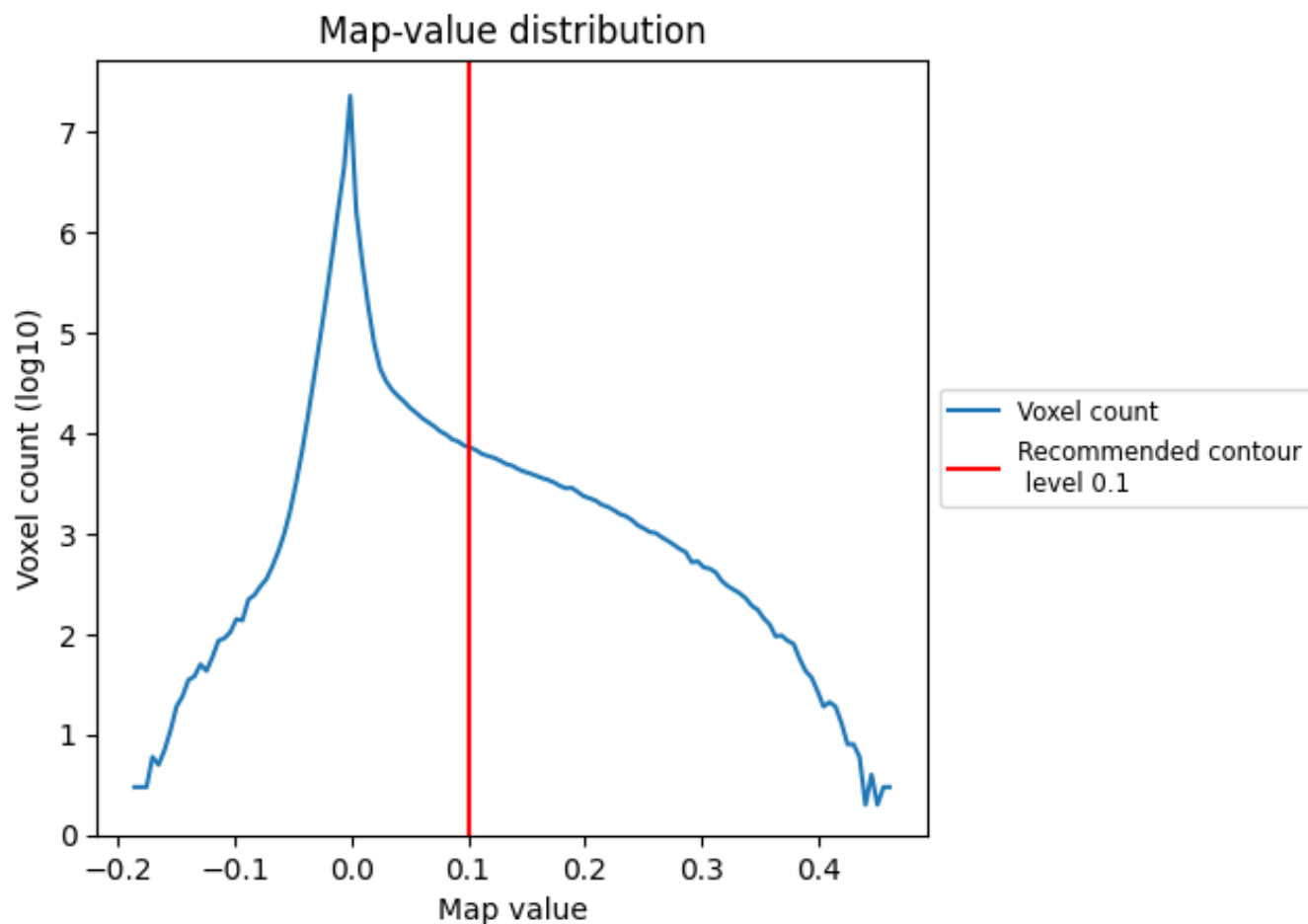
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

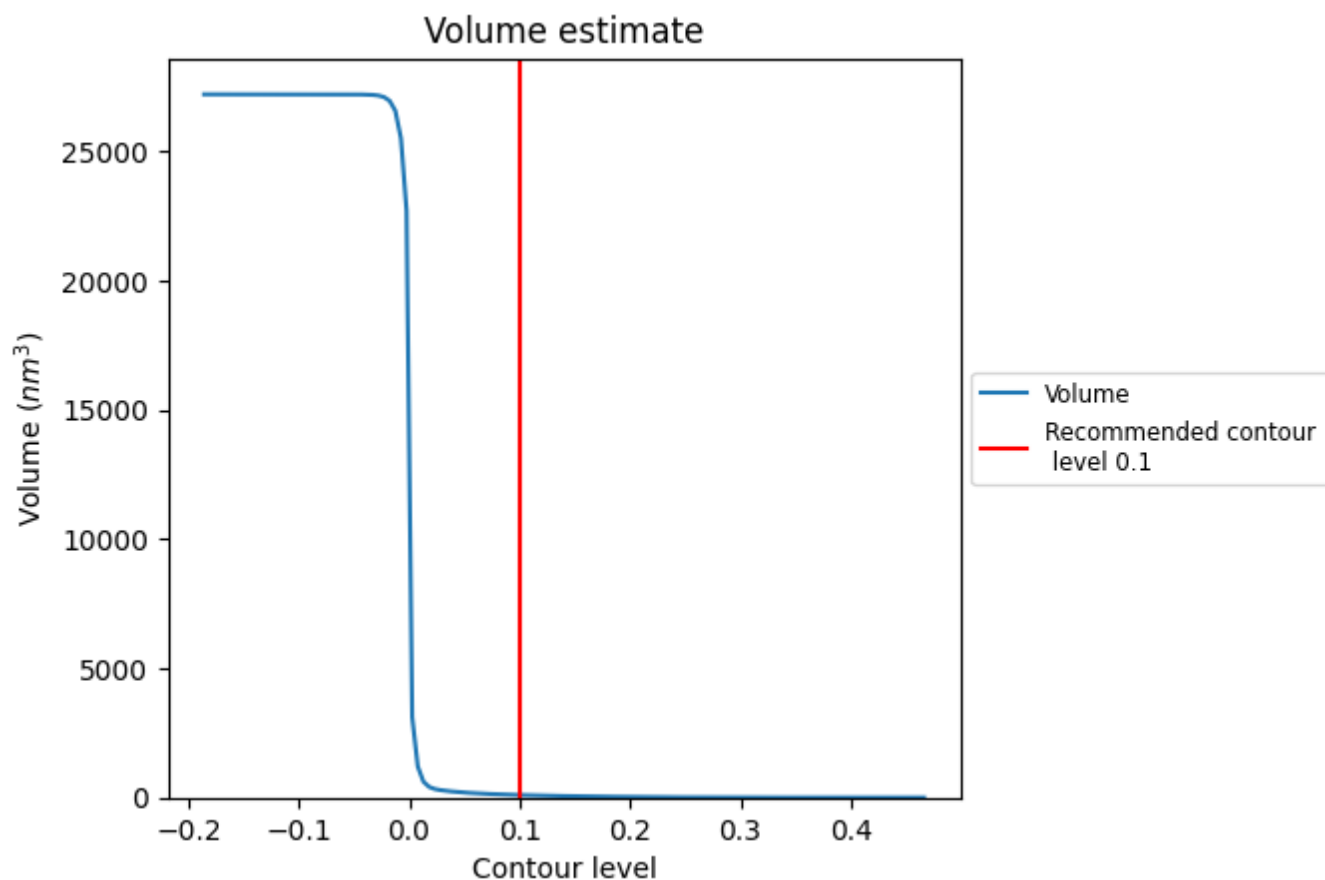
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

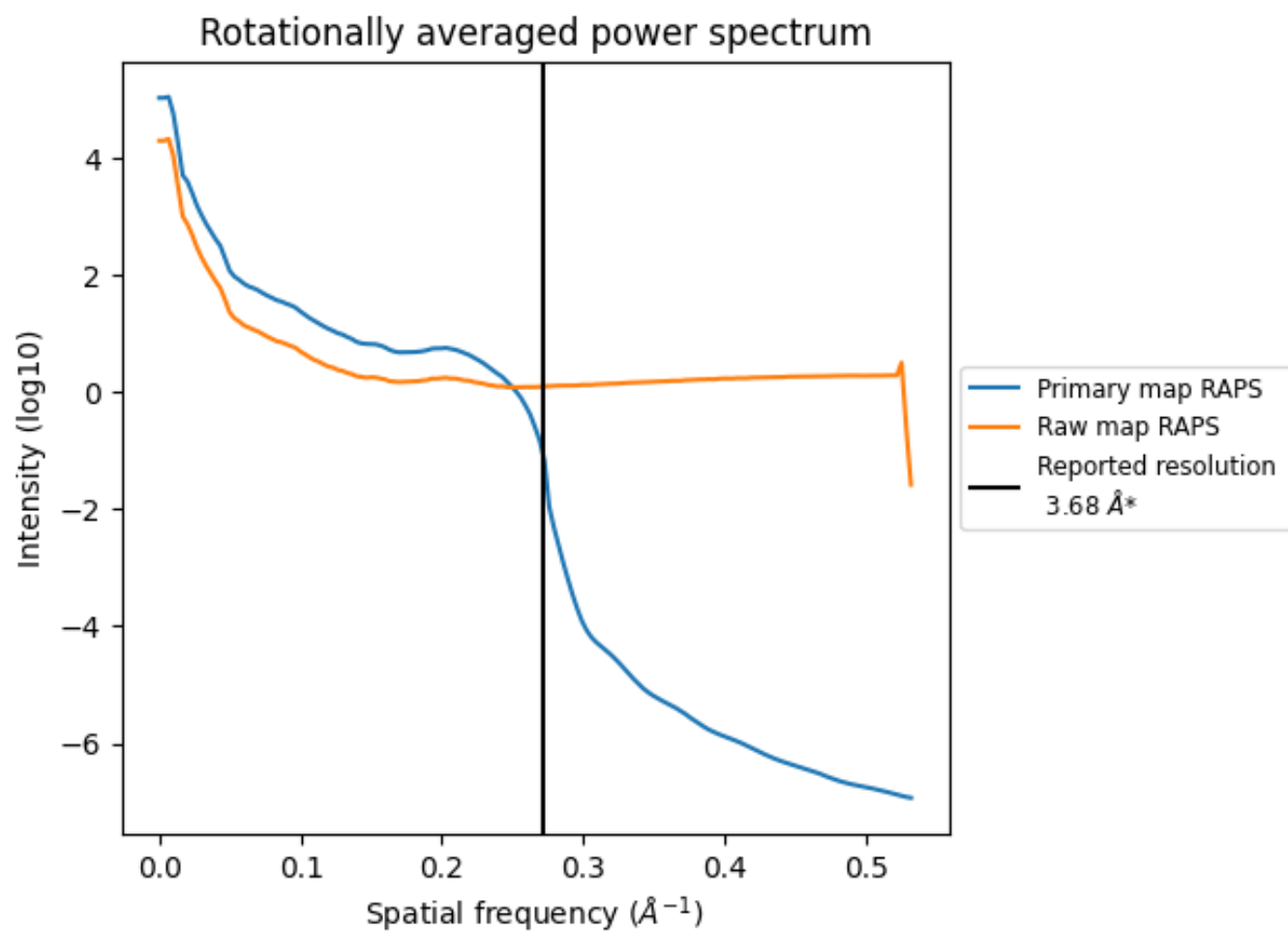
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 97 nm^3 ; this corresponds to an approximate mass of 88 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 ⓘ

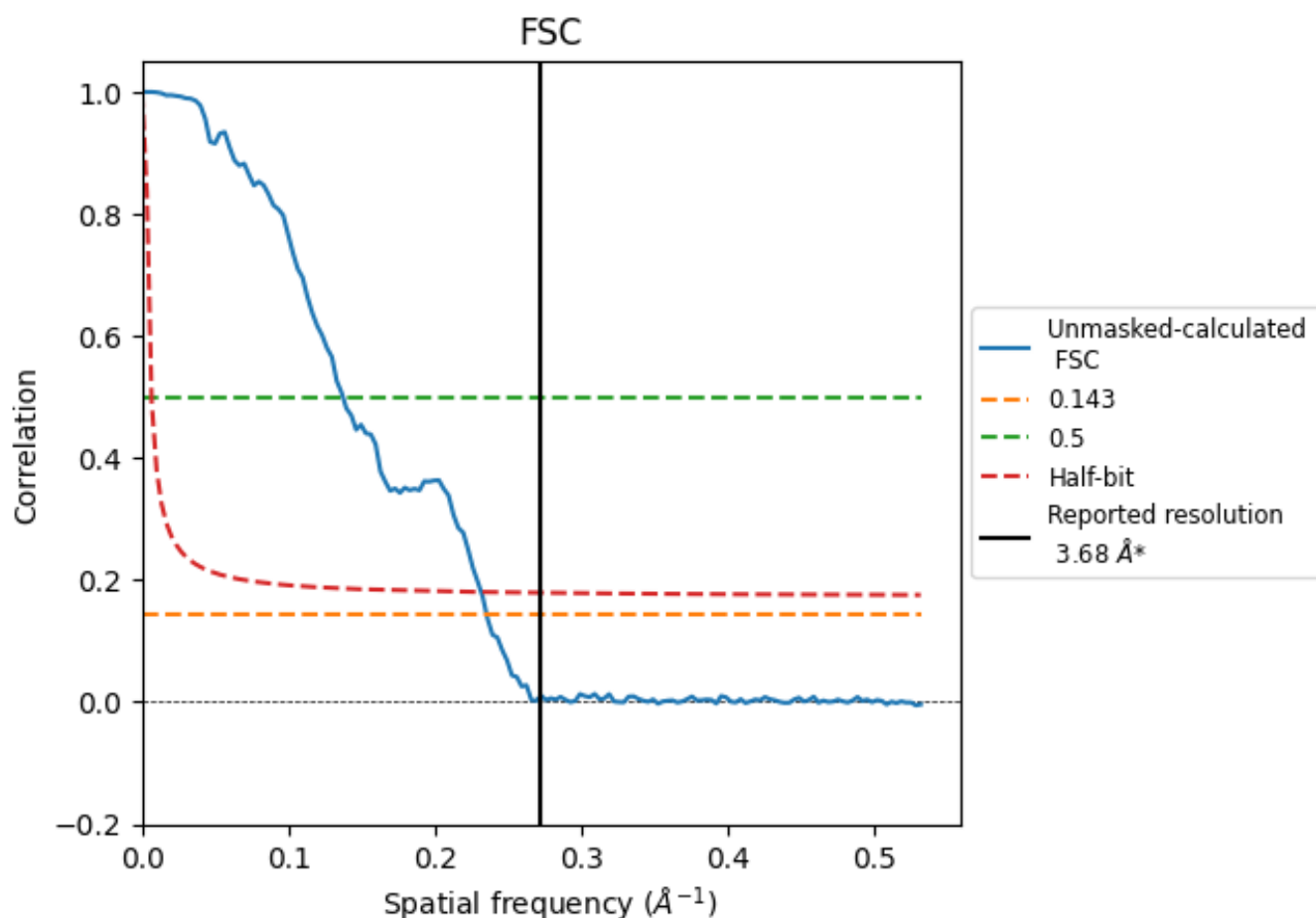


*Reported resolution corresponds to spatial frequency of 0.272 \AA^{-1}

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 0.272 \AA^{-1}

8.2 Resolution estimates [i](#)

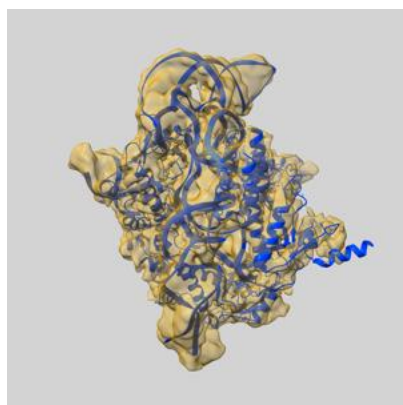
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.68	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.25	7.29	4.32

*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 4.25 differs from the reported value 3.68 by more than 10 %

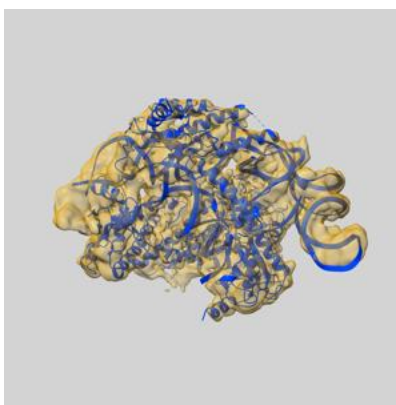
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-49958 and PDB model 9NZS. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

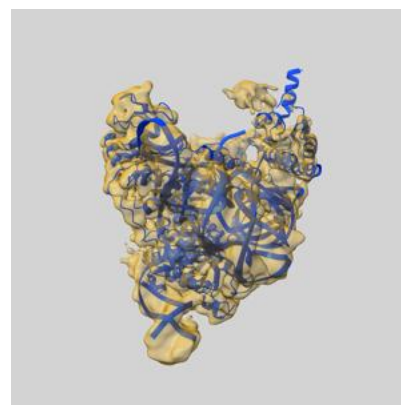
9.1 Map-model overlay [i](#)



X



Y



Z

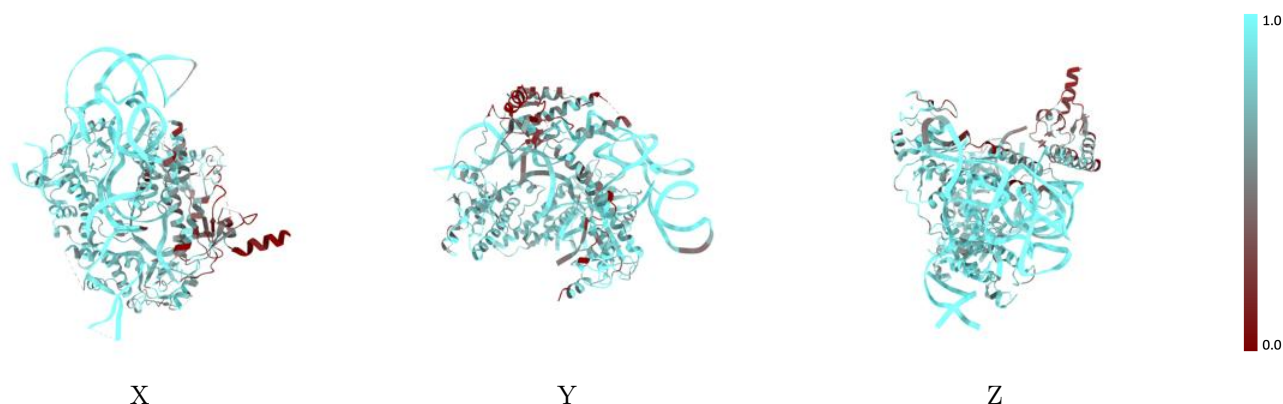
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



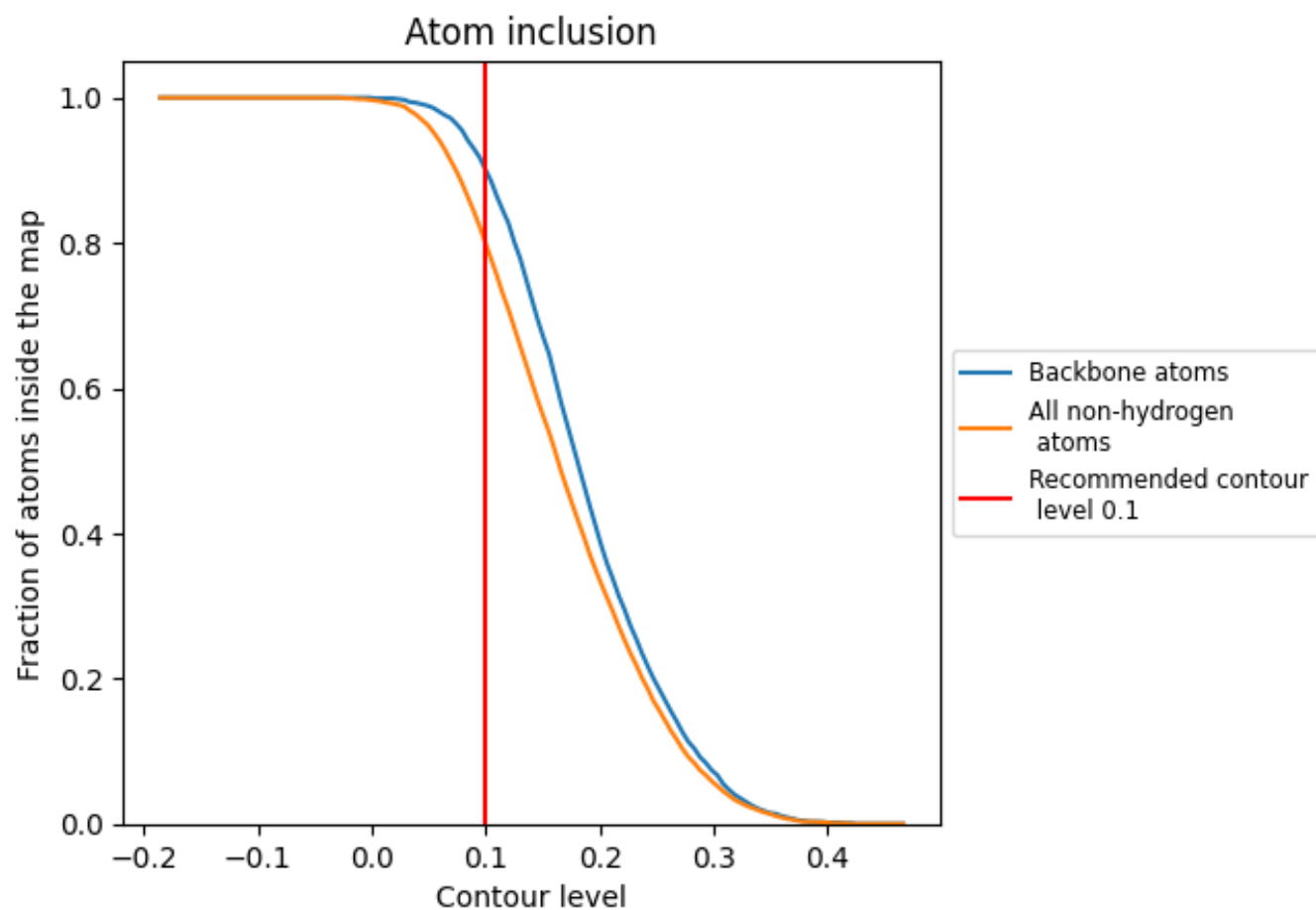
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 80% 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 (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div>0.7980</div>	<div><div></div>0.3510</div>
A	<div><div></div>0.8040</div>	<div><div></div>0.3970</div>
B	<div><div></div>0.6280</div>	<div><div></div>0.2940</div>
C	<div><div></div>0.8540</div>	<div><div></div>0.3610</div>
D	<div><div></div>0.9820</div>	<div><div></div>0.3930</div>
E	<div><div></div>0.9290</div>	<div><div></div>0.3460</div>
G	<div><div></div>0.4210</div>	<div><div></div>0.2880</div>

1.0

0.0

<0.0