



Full wwPDB EM Validation Report ⓘ

Apr 15, 2026 – 12:04 PM UTC

PDB ID : 7JZV / pdb_00007jzv
EMDB ID : EMD-22581
Title : Cryo-EM structure of the BRCA1-UbcH5c/BARD1 E3-E2 module bound to a nucleosome
Authors : Witus, S.R.; Burrell, A.L.; Hansen, J.M.; Farrell, D.P.; Dimaio, F.; Kollman, J.M.; Klevit, R.E.
Deposited on : 2020-09-02
Resolution : 3.90 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:

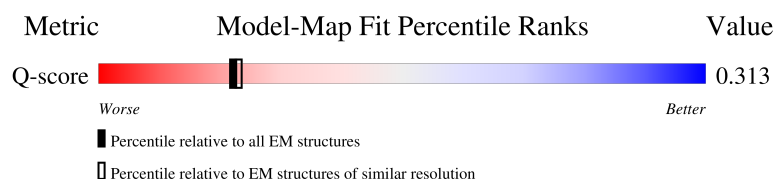
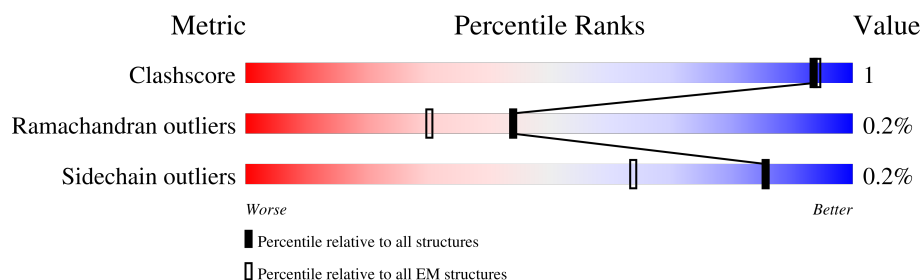
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.90 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.



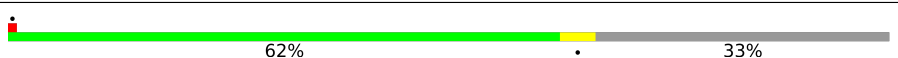
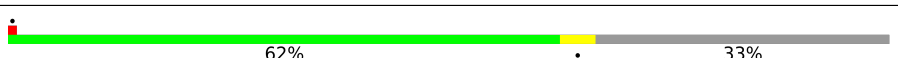
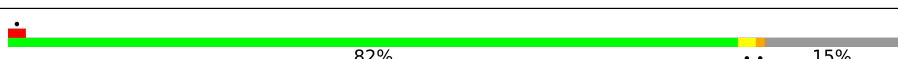
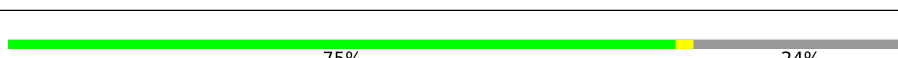
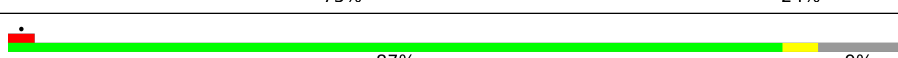
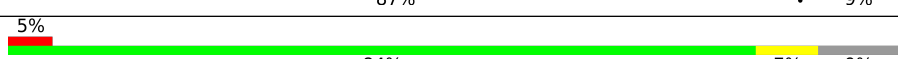
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	8855 (3.40 - 4.40)

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	258	40% 87% 10% .
2	B	115	20% 81% 16% ..
3	N	133	71% 8% 22%
3	n	133	5% 72% 6% 22%

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Mol	Chain	Length	Quality of chain
4	O	125	
4	o	125	
5	P	135	
5	p	135	
6	Q	102	
6	q	102	
7	X	153	
8	Y	153	

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14339 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 BRCA1,Ubiquitin-conjugating enzyme E2 D3.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	253	Total	C	N	O	S	0	0
			1995	1268	340	371	16		

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP A0A386IN42
A	1	PRO	-	expression tag	UNP A0A386IN42
A	105	GLY	-	linker	UNP A0A386IN42
A	106	SER	-	linker	UNP A0A386IN42
A	107	GLY	-	linker	UNP A0A386IN42
A	108	SER	-	linker	UNP A0A386IN42
A	109	GLY	-	linker	UNP A0A386IN42
A	110	SER	-	linker	UNP A0A386IN42
A	111	GLY	-	linker	UNP A0A386IN42
A	285	LYS	CYS	engineered mutation	UNP P61077

- Molecule 2 is a protein called BRCA1-associated RING domain protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	97	Total	C	N	O	S	0	0
			753	461	140	140	12		

- Molecule 3 is a protein called Histone H2A type 2-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	N	104	Total	C	N	O	S	0	0
			801	506	155	139	1		
3	n	104	Total	C	N	O	S	0	0
			801	506	155	139	1		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
N	-3	GLY	-	expression tag	UNP Q6FI13
N	-2	ALA	-	expression tag	UNP Q6FI13
N	-1	MET	-	expression tag	UNP Q6FI13
N	0	GLY	-	expression tag	UNP Q6FI13
n	-3	GLY	-	expression tag	UNP Q6FI13
n	-2	ALA	-	expression tag	UNP Q6FI13
n	-1	MET	-	expression tag	UNP Q6FI13
n	0	GLY	-	expression tag	UNP Q6FI13

- Molecule 4 is a protein called Histone H2B type 1-K.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	O	94	Total	C	N	O	S	0	0
			735	461	134	138	2		
4	o	96	Total	C	N	O	S	0	0
			754	473	138	141	2		

- Molecule 5 is a protein called Histone H3.2.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	P	90	Total	C	N	O	S	0	0
			737	465	140	130	2		
5	p	90	Total	C	N	O	S	0	0
			738	466	140	130	2		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	110	ALA	CYS	engineered mutation	UNP Q71DI3
p	110	ALA	CYS	engineered mutation	UNP Q71DI3

- Molecule 6 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	Q	87	Total	C	N	O	S	0	0
			703	442	142	118	1		
6	q	78	Total	C	N	O	S	0	0
			619	391	120	107	1		

- Molecule 7 is a DNA chain called Widom 601 153-bp.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	X	139	Total	C	N	O	P	0	0
			2867	1357	536	835	139		

- Molecule 8 is a DNA chain called Widom 601 153-bp.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	Y	139	Total	C	N	O	P	0	0
			2832	1345	515	833	139		

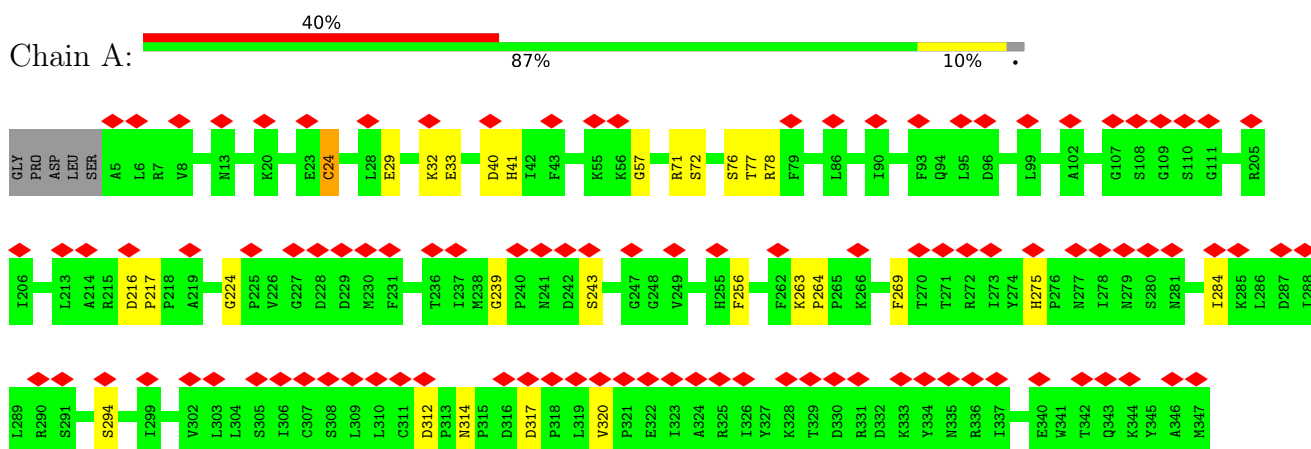
- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
9	A	2	Total	Zn	0
			2	2	
9	B	2	Total	Zn	0
			2	2	

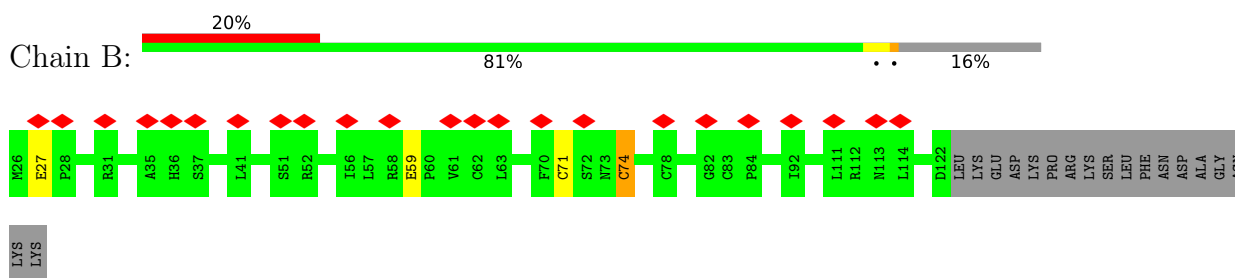
3 Residue-property plots

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.

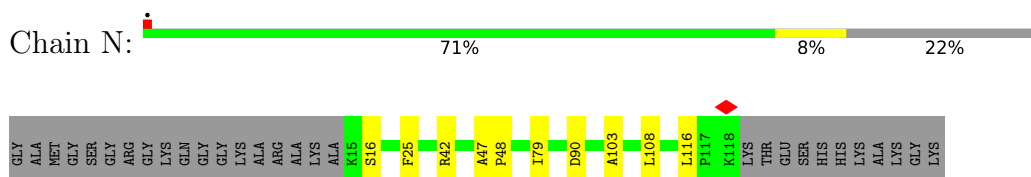
- Molecule 1: BRCA1,Ubiquitin-conjugating enzyme E2 D3



- Molecule 2: BRCA1-associated RING domain protein 1

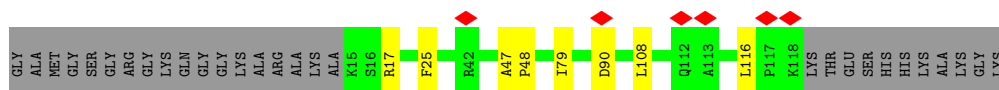


- Molecule 3: Histone H2A type 2-A

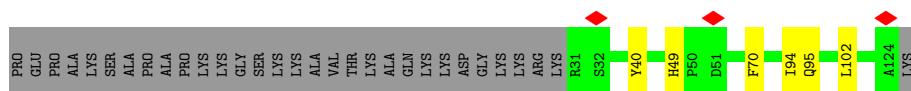


- Molecule 3: Histone H2A type 2-A

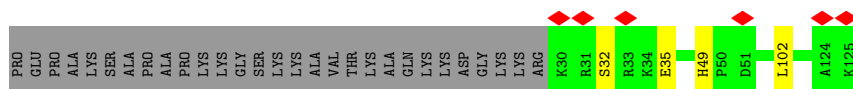




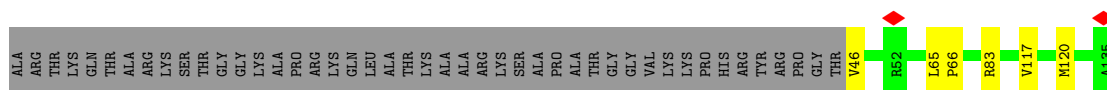
• Molecule 4: Histone H2B type 1-K



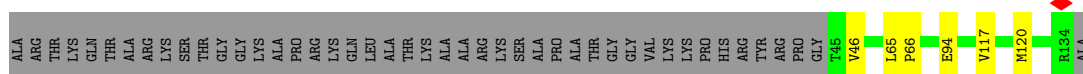
• Molecule 4: Histone H2B type 1-K



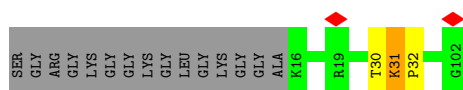
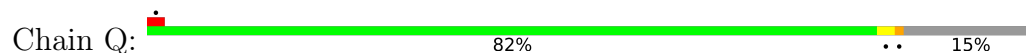
• Molecule 5: Histone H3.2



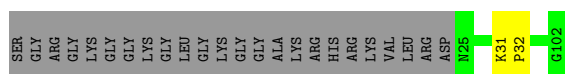
• Molecule 5: Histone H3.2



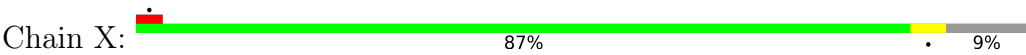
• Molecule 6: Histone H4



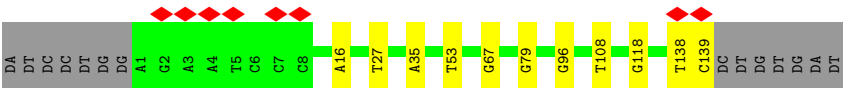
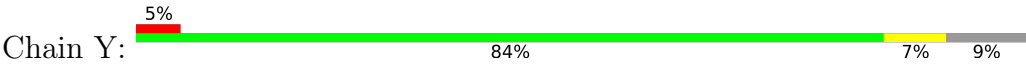
• Molecule 6: Histone H4



• Molecule 7: Widom 601 153-bp



• Molecule 8: Widom 601 153-bp



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	21479	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	90	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2700	Depositor
Magnification	130000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	14.926	Depositor
Minimum map value	-11.777	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.341	Depositor
Recommended contour level	0.7	Depositor
Map size (Å)	268.8, 268.8, 268.8	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.05, 1.05, 1.05	Depositor

5 Model quality

5.1 Standard geometry

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	0.66	0/2040	1.09	35/2763 (1.3%)
2	B	0.85	0/765	1.04	5/1033 (0.5%)
3	N	0.83	0/811	0.91	8/1093 (0.7%)
3	n	0.77	0/811	0.86	8/1093 (0.7%)
4	O	0.88	0/746	0.89	4/1003 (0.4%)
4	o	0.77	0/765	0.80	4/1025 (0.4%)
5	P	0.86	0/745	0.74	2/997 (0.2%)
5	p	0.88	0/746	0.77	2/1000 (0.2%)
6	Q	0.84	0/711	0.78	2/948 (0.2%)
6	q	0.85	0/626	0.77	0/837
7	X	0.45	0/3219	0.94	0/4971
8	Y	0.46	0/3173	0.94	3/4890 (0.1%)
All	All	0.68	0/15158	0.92	73/21653 (0.3%)

There are no bond length outliers.

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	263	LYS	CA-C-N	8.69	126.00	119.66
1	A	263	LYS	C-N-CA	8.69	126.00	119.66
4	O	49	HIS	CA-C-N	8.56	128.20	119.56
4	O	49	HIS	C-N-CA	8.56	128.20	119.56
1	A	264	PRO	CA-C-N	7.91	128.06	120.31
1	A	264	PRO	C-N-CA	7.91	128.06	120.31
1	A	312	ASP	CA-C-N	7.73	128.19	119.92
1	A	312	ASP	C-N-CA	7.73	128.19	119.92
1	A	239	GLY	CA-C-N	7.37	127.81	119.92
1	A	239	GLY	C-N-CA	7.37	127.81	119.92
4	o	49	HIS	CA-C-N	7.32	127.33	119.28
4	o	49	HIS	C-N-CA	7.32	127.33	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	317	ASP	CA-C-N	7.26	127.71	120.52
1	A	317	ASP	C-N-CA	7.26	127.71	120.52
1	A	216	ASP	CA-C-N	6.96	127.55	120.38
1	A	216	ASP	C-N-CA	6.96	127.55	120.38
4	O	102	LEU	CA-C-N	6.93	126.89	120.03
4	O	102	LEU	C-N-CA	6.93	126.89	120.03
3	n	116	LEU	CA-C-N	6.82	126.80	119.78
3	n	116	LEU	C-N-CA	6.82	126.80	119.78
2	B	27	GLU	CA-C-N	6.78	127.04	119.32
2	B	27	GLU	C-N-CA	6.78	127.04	119.32
3	N	116	LEU	CA-C-N	6.71	126.69	119.78
3	N	116	LEU	C-N-CA	6.71	126.69	119.78
1	A	275	HIS	CA-C-N	6.65	127.19	119.47
1	A	275	HIS	C-N-CA	6.65	127.19	119.47
2	B	59	GLU	CA-C-N	6.56	126.32	119.76
2	B	59	GLU	C-N-CA	6.56	126.32	119.76
1	A	256	PHE	CA-C-N	6.48	126.66	120.31
1	A	256	PHE	C-N-CA	6.48	126.66	120.31
1	A	320	VAL	CA-C-N	6.41	126.91	119.47
1	A	320	VAL	C-N-CA	6.41	126.91	119.47
1	A	224	GLY	CA-C-N	6.40	126.32	119.85
1	A	224	GLY	C-N-CA	6.40	126.32	119.85
1	A	24	CYS	CA-C-N	6.30	126.78	119.47
1	A	24	CYS	C-N-CA	6.30	126.78	119.47
1	A	33	GLU	CA-C-N	6.29	126.05	119.76
1	A	33	GLU	C-N-CA	6.29	126.05	119.76
1	A	217	PRO	CA-C-N	6.27	126.30	119.90
1	A	217	PRO	C-N-CA	6.27	126.30	119.90
3	N	25	PHE	CA-C-N	5.90	127.22	119.84
3	N	25	PHE	C-N-CA	5.90	127.22	119.84
1	A	243	SER	CA-C-N	5.79	125.89	119.87
1	A	243	SER	C-N-CA	5.79	125.89	119.87
1	A	57	GLY	CA-C-N	5.75	126.08	119.93
1	A	57	GLY	C-N-CA	5.75	126.08	119.93
3	n	25	PHE	CA-C-N	5.74	127.02	119.84
3	n	25	PHE	C-N-CA	5.74	127.02	119.84
4	o	102	LEU	CA-C-N	5.74	126.22	120.14
4	o	102	LEU	C-N-CA	5.74	126.22	120.14
1	A	294	SER	CA-C-N	5.64	126.89	119.84
1	A	294	SER	C-N-CA	5.64	126.89	119.84
1	A	314	ASN	CA-C-N	5.55	126.61	120.45
1	A	314	ASN	C-N-CA	5.55	126.61	120.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	n	79	ILE	CA-C-N	5.46	125.55	119.32
3	n	79	ILE	C-N-CA	5.46	125.55	119.32
5	p	120	MET	CA-C-N	5.39	125.47	119.32
5	p	120	MET	C-N-CA	5.39	125.47	119.32
3	N	108	LEU	CA-C-N	5.33	125.63	119.93
3	N	108	LEU	C-N-CA	5.33	125.63	119.93
2	B	74	CYS	CA-CB-SG	5.26	126.50	114.40
3	N	79	ILE	CA-C-N	5.21	125.26	119.32
3	N	79	ILE	C-N-CA	5.21	125.26	119.32
6	Q	31	LYS	CA-C-N	5.17	125.22	119.32
6	Q	31	LYS	C-N-CA	5.17	125.22	119.32
5	P	120	MET	CA-C-N	5.11	125.39	119.47
5	P	120	MET	C-N-CA	5.11	125.39	119.47
1	A	40	ASP	N-CA-C	5.10	117.50	109.39
8	Y	53	DT	P-O3'-C3'	5.07	127.81	120.20
8	Y	35	DA	P-O3'-C3'	5.07	127.80	120.20
3	n	108	LEU	CA-C-N	5.04	125.32	119.93
3	n	108	LEU	C-N-CA	5.04	125.32	119.93
8	Y	16	DA	P-O3'-C3'	5.03	127.74	120.20

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	1995	0	2007	5	0
2	B	753	0	737	3	0
3	N	801	0	857	5	0
3	n	801	0	857	3	0
4	O	735	0	756	3	0
4	o	754	0	782	2	0
5	P	737	0	777	4	0
5	p	738	0	779	4	0
6	Q	703	0	755	5	0
6	q	619	0	659	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	X	2867	0	1561	7	0
8	Y	2832	0	1560	7	0
9	A	2	0	0	0	0
9	B	2	0	0	0	0
All	All	14339	0	12087	35	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 1.

All (35) 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)
8:Y:27:DT:OP2	3:n:17:ARG:HD3	1.85	0.76
7:X:120:DG:OP1	4:o:32:SER:HB2	1.88	0.74
6:Q:30:THR:OG1	7:X:58:DA:OP2	2.12	0.68
5:P:46:VAL:HB	8:Y:79:DG:OP1	1.93	0.68
7:X:120:DG:OP1	4:o:32:SER:CB	2.46	0.64
3:N:16:SER:HA	7:X:27:DA:H5''	1.84	0.58
8:Y:67:DG:OP1	5:p:117:VAL:N	2.31	0.58
6:Q:31:LYS:HB2	6:Q:32:PRO:HD3	1.87	0.56
6:q:31:LYS:HB3	6:q:32:PRO:HD3	1.87	0.56
2:B:71:CYS:HB2	2:B:74:CYS:HB2	1.89	0.54
3:n:90:ASP:OD1	3:n:90:ASP:C	2.53	0.51
4:O:40:TYR:OH	8:Y:118:DG:H5''	2.11	0.50
3:N:42:ARG:HH21	8:Y:108:DT:C1'	2.26	0.48
1:A:71:ARG:O	1:A:72:SER:OG	2.30	0.48
3:n:47:ALA:N	3:n:48:PRO:HD2	2.28	0.48
3:N:47:ALA:N	3:N:48:PRO:HD2	2.29	0.48
1:A:71:ARG:O	1:A:72:SER:CB	2.61	0.47
6:Q:30:THR:OG1	6:Q:31:LYS:N	2.47	0.47
6:Q:30:THR:HG21	7:X:57:DA:H3'	1.97	0.47
5:P:65:LEU:N	5:P:66:PRO:HD2	2.30	0.47
1:A:269:PHE:CD1	1:A:284:ILE:HD11	2.50	0.46
5:P:83:ARG:HH22	8:Y:96:DG:C4'	2.29	0.46
8:Y:138:DT:H2''	8:Y:139:DC:H5'	1.98	0.46
5:P:117:VAL:N	7:X:67:DA:OP1	2.44	0.45
5:p:65:LEU:N	5:p:66:PRO:HD2	2.32	0.43
2:B:71:CYS:CB	2:B:74:CYS:HB2	2.49	0.43
1:A:76:SER:O	1:A:78:ARG:N	2.52	0.42
4:O:94:ILE:O	4:O:95:GLN:C	2.62	0.42
7:X:79:DT:OP1	5:p:46:VAL:HB	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:Q:31:LYS:CB	6:Q:32:PRO:HD3	2.51	0.41
2:B:71:CYS:N	2:B:74:CYS:HB3	2.35	0.40
3:N:103:ALA:HA	5:p:94:GLU:CD	2.47	0.40
4:O:70:PHE:CD2	4:O:70:PHE:C	2.99	0.40
1:A:29:GLU:OE1	1:A:32:LYS:NZ	2.55	0.40
3:N:90:ASP:OD1	3:N:90:ASP:C	2.65	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	251/258 (97%)	245 (98%)	5 (2%)	1 (0%)	30	64
2	B	95/115 (83%)	89 (94%)	6 (6%)	0	100	100
3	N	102/133 (77%)	100 (98%)	2 (2%)	0	100	100
3	n	102/133 (77%)	100 (98%)	2 (2%)	0	100	100
4	O	92/125 (74%)	92 (100%)	0	0	100	100
4	o	94/125 (75%)	93 (99%)	0	1 (1%)	11	43
5	P	88/135 (65%)	87 (99%)	1 (1%)	0	100	100
5	p	88/135 (65%)	88 (100%)	0	0	100	100
6	Q	85/102 (83%)	85 (100%)	0	0	100	100
6	q	76/102 (74%)	76 (100%)	0	0	100	100
All	All	1073/1363 (79%)	1055 (98%)	16 (2%)	2 (0%)	44	74

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	77	THR

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Mol	Chain	Res	Type
4	o	35	GLU

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	228/232 (98%)	226 (99%)	2 (1%)	70	75
2	B	86/102 (84%)	86 (100%)	0	100	100
3	N	82/99 (83%)	82 (100%)	0	100	100
3	n	82/99 (83%)	82 (100%)	0	100	100
4	O	80/104 (77%)	80 (100%)	0	100	100
4	o	82/104 (79%)	82 (100%)	0	100	100
5	P	77/109 (71%)	77 (100%)	0	100	100
5	p	78/109 (72%)	78 (100%)	0	100	100
6	Q	72/78 (92%)	72 (100%)	0	100	100
6	q	63/78 (81%)	63 (100%)	0	100	100
All	All	930/1114 (84%)	928 (100%)	2 (0%)	85	87

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

Mol	Chain	Res	Type
1	A	24	CYS
1	A	41	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
3	N	24	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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 4 ligands modelled in this entry, 4 are 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.

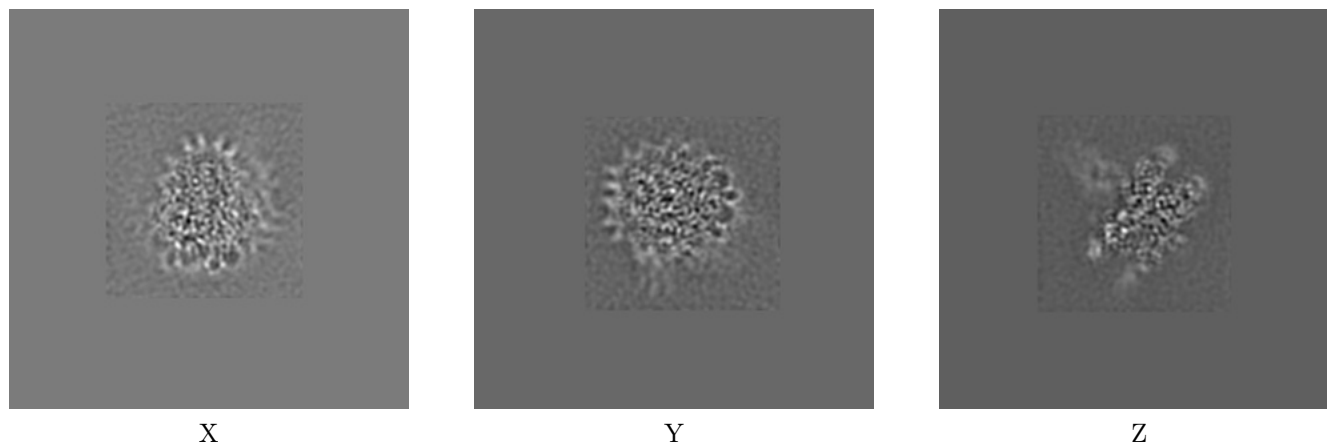
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-22581. These allow visual inspection of the internal detail of the map and identification of artifacts.

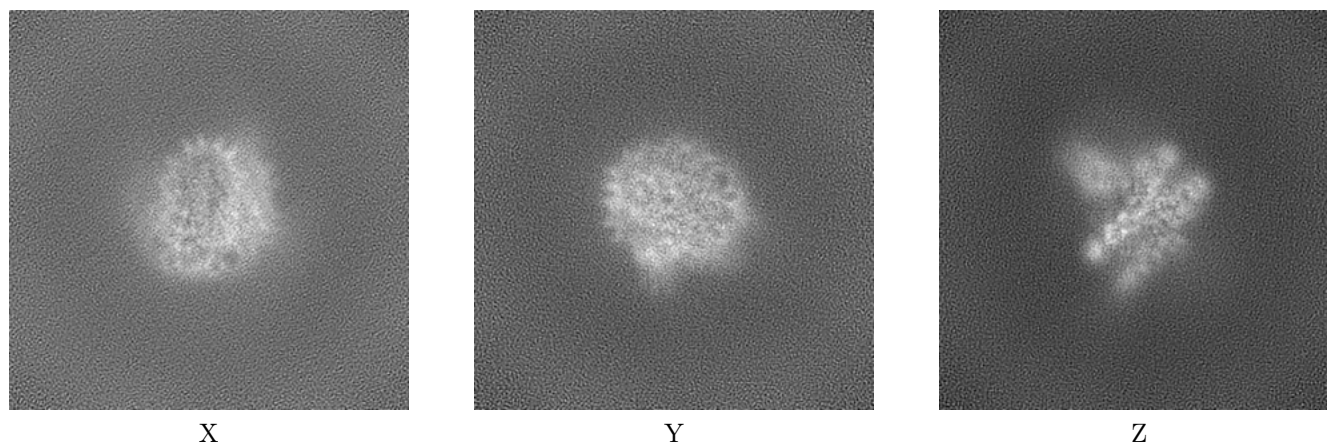
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

6.1.1 Primary map



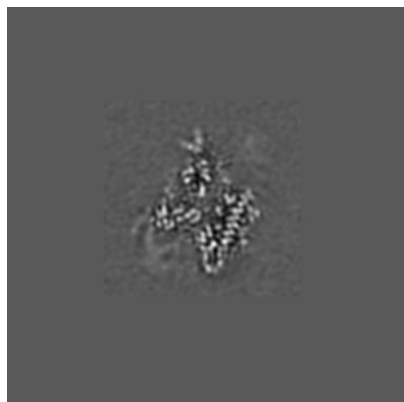
6.1.2 Raw map



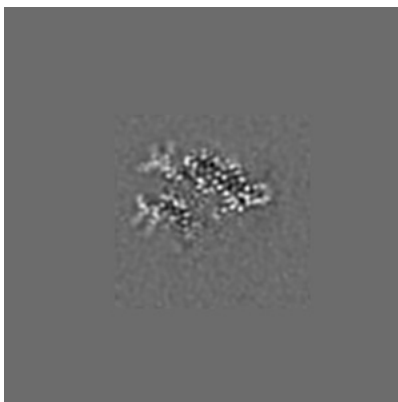
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

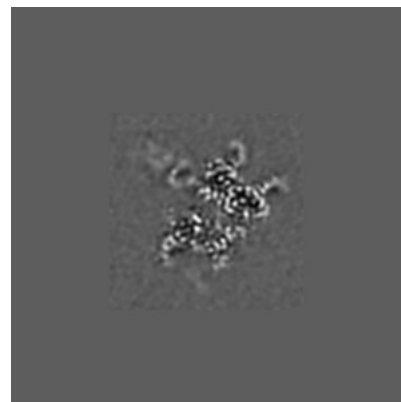
6.2.1 Primary map



X Index: 128

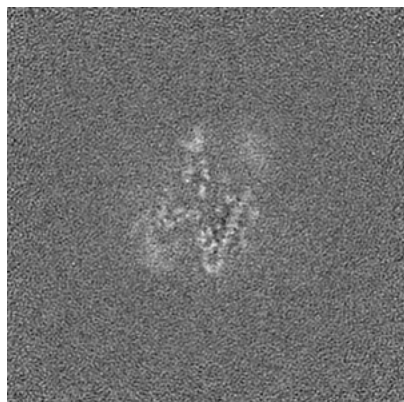


Y Index: 128

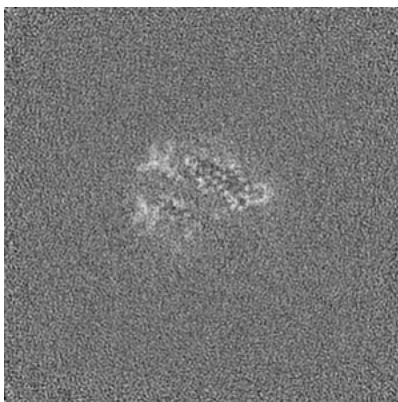


Z Index: 128

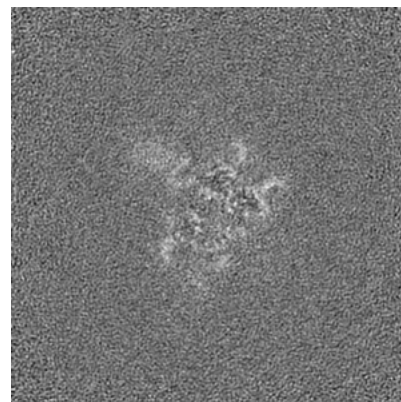
6.2.2 Raw map



X Index: 128



Y Index: 128

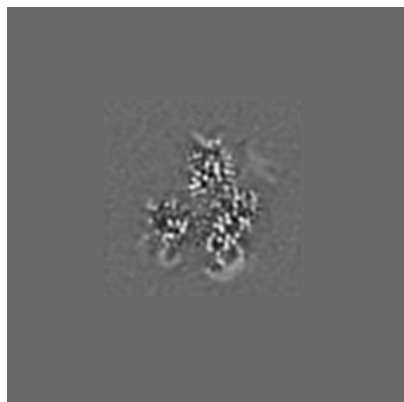


Z Index: 128

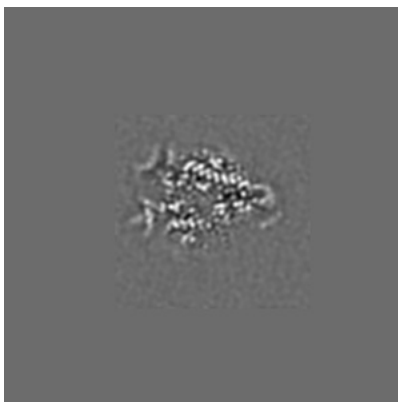
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

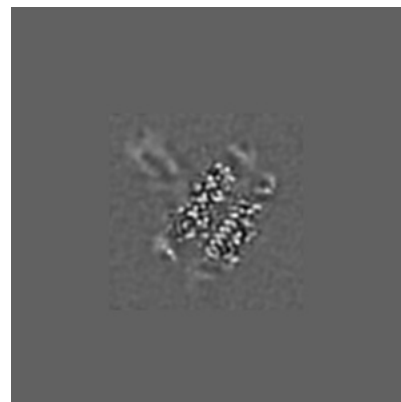
6.3.1 Primary map



X Index: 134

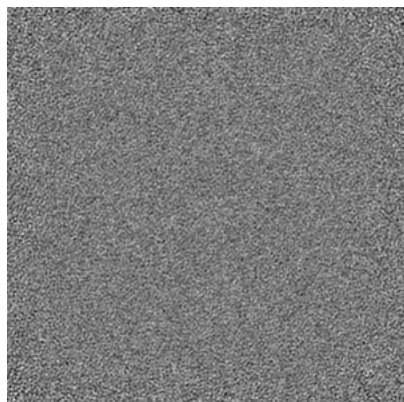


Y Index: 125

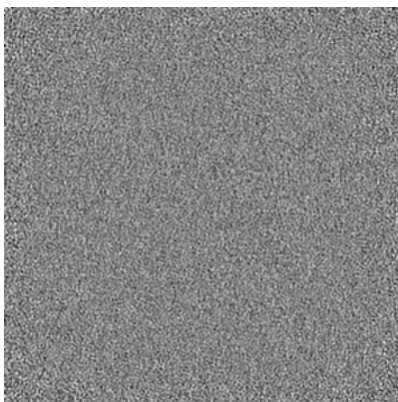


Z Index: 116

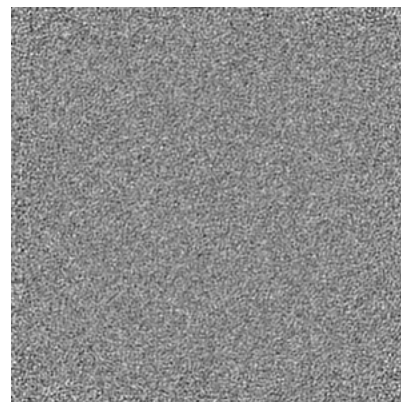
6.3.2 Raw map



X Index: 0



Y Index: 0

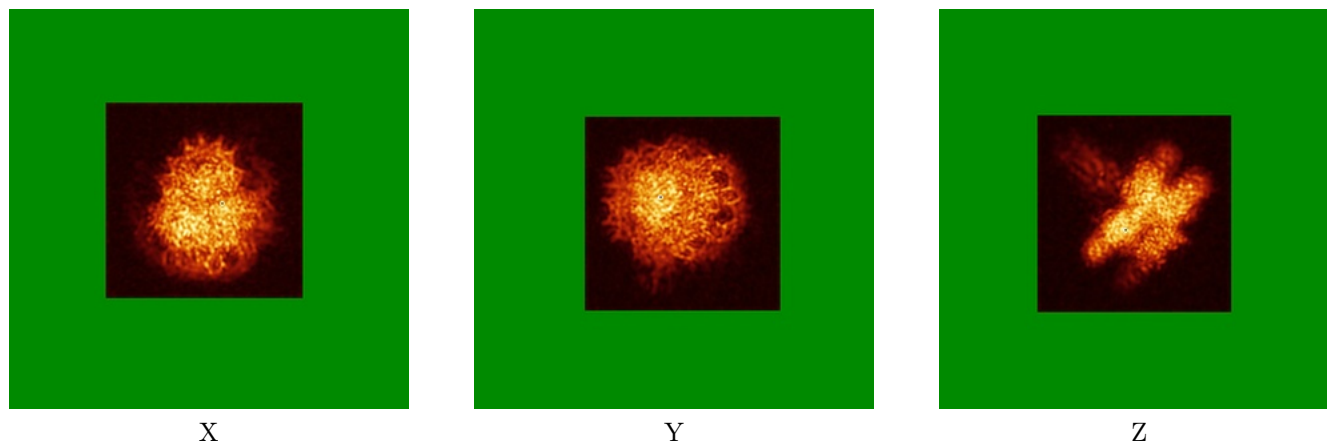


Z Index: 0

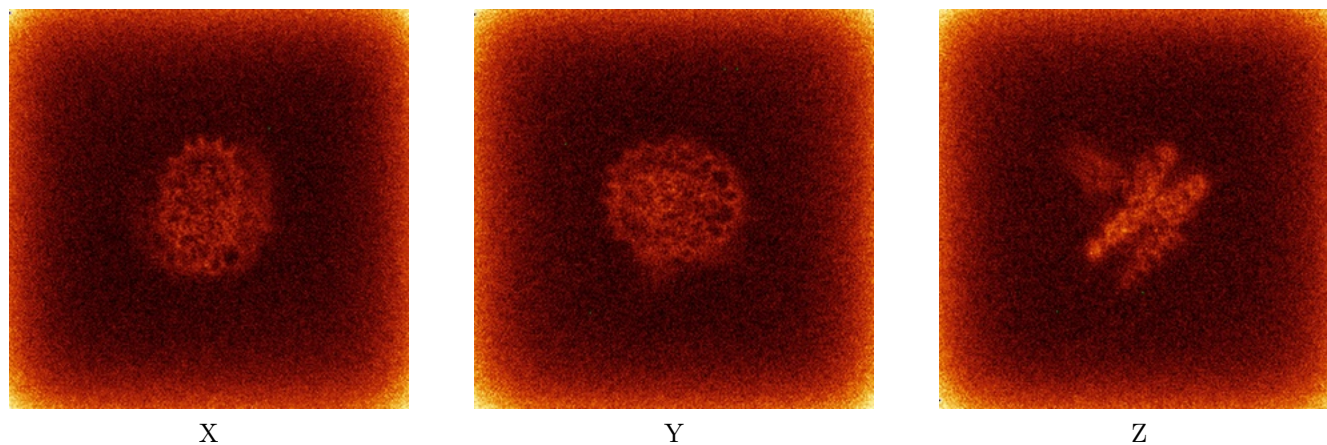
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



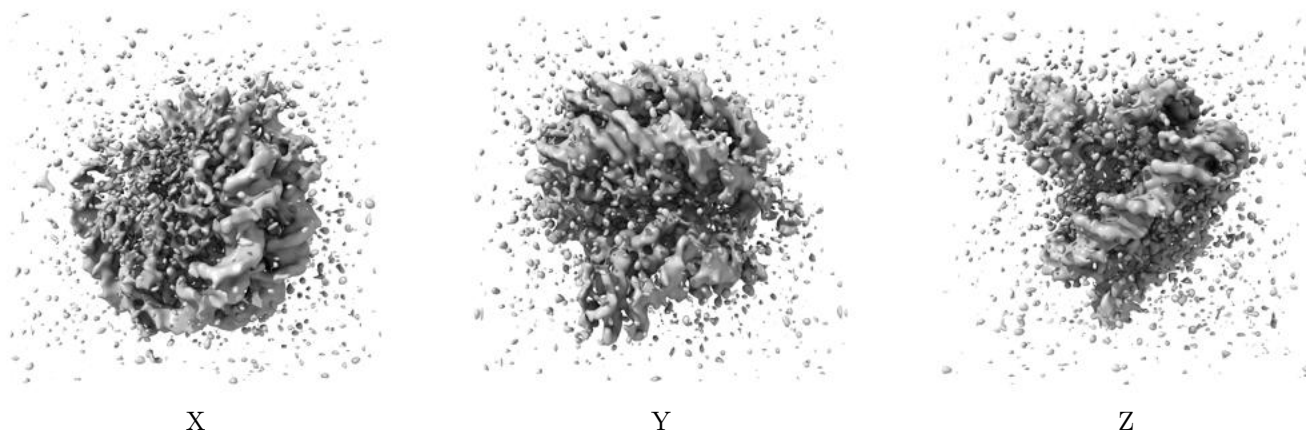
6.4.2 Raw map



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.

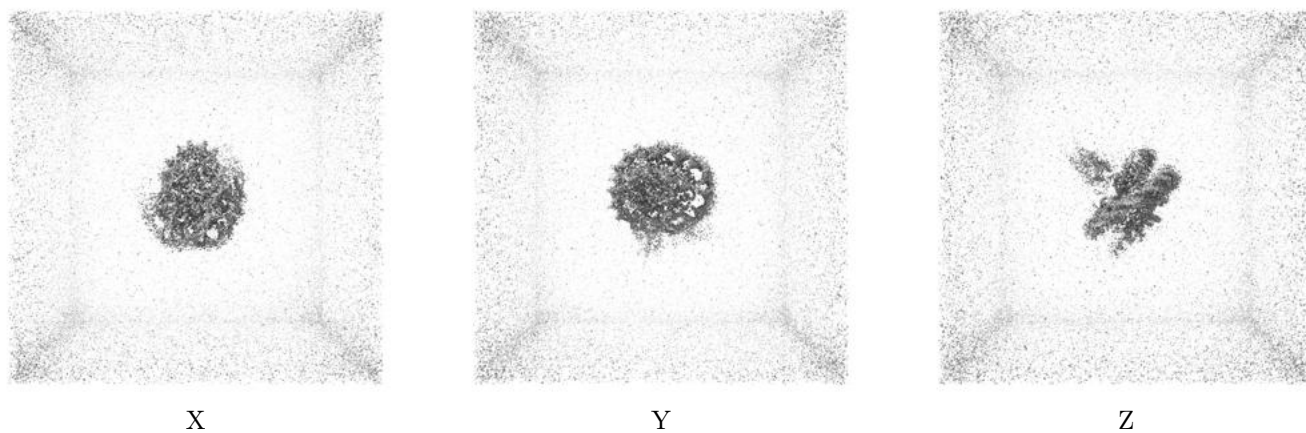
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.7. 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.

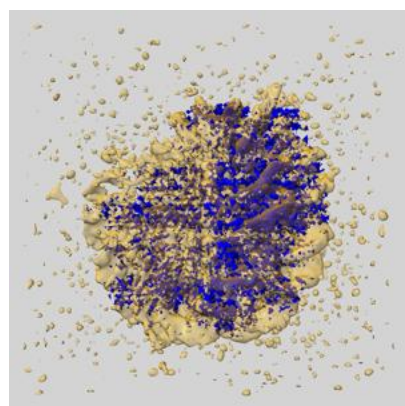
6.6 Mask visualisation [i](#)

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

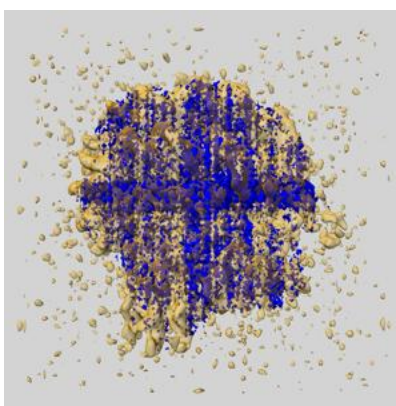
A mask typically either:

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

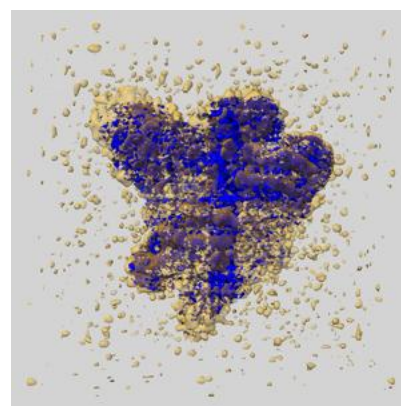
6.6.1 emd_22581_msk_1.map [i](#)



X



Y

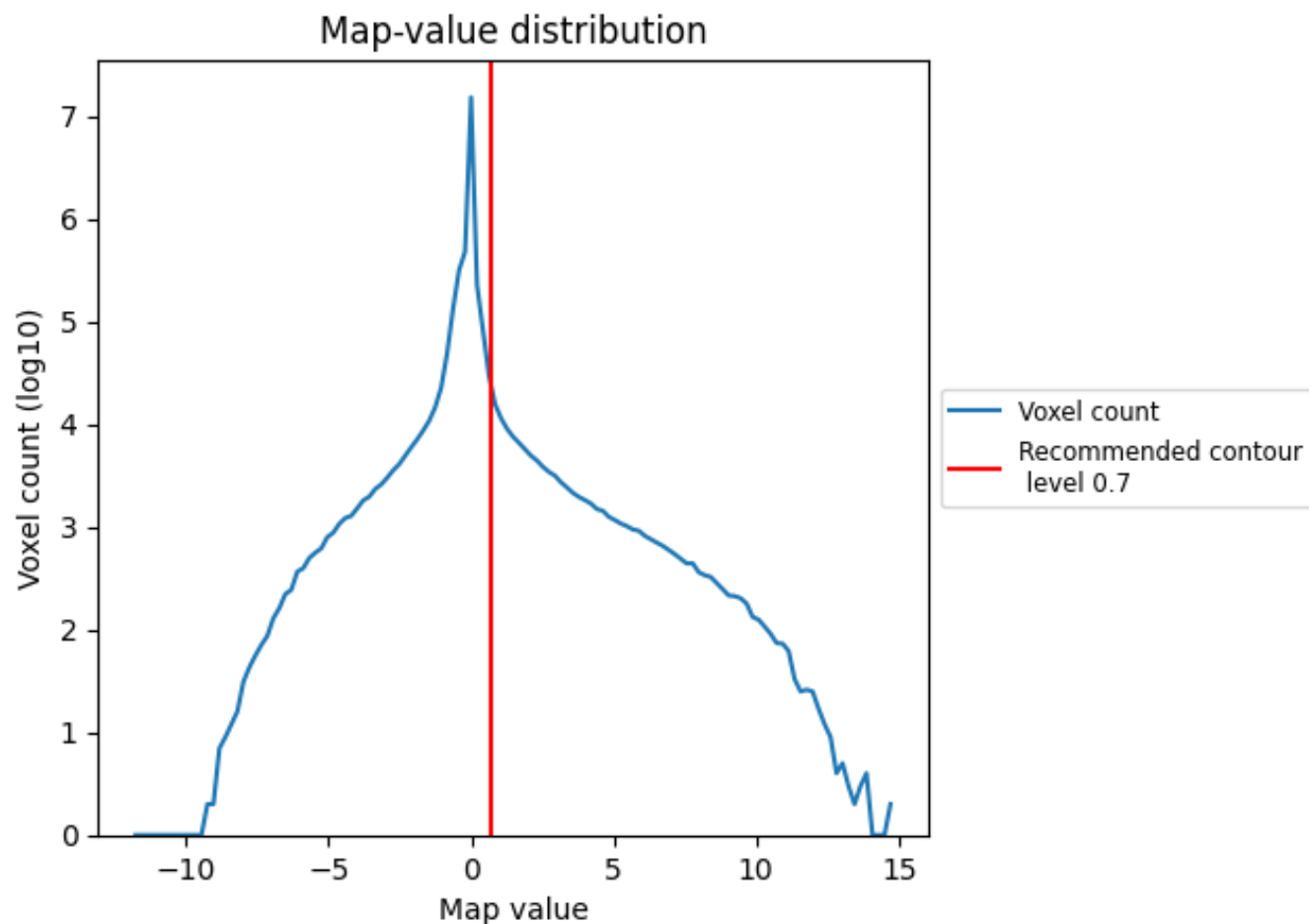


Z

7 Map analysis [i](#)

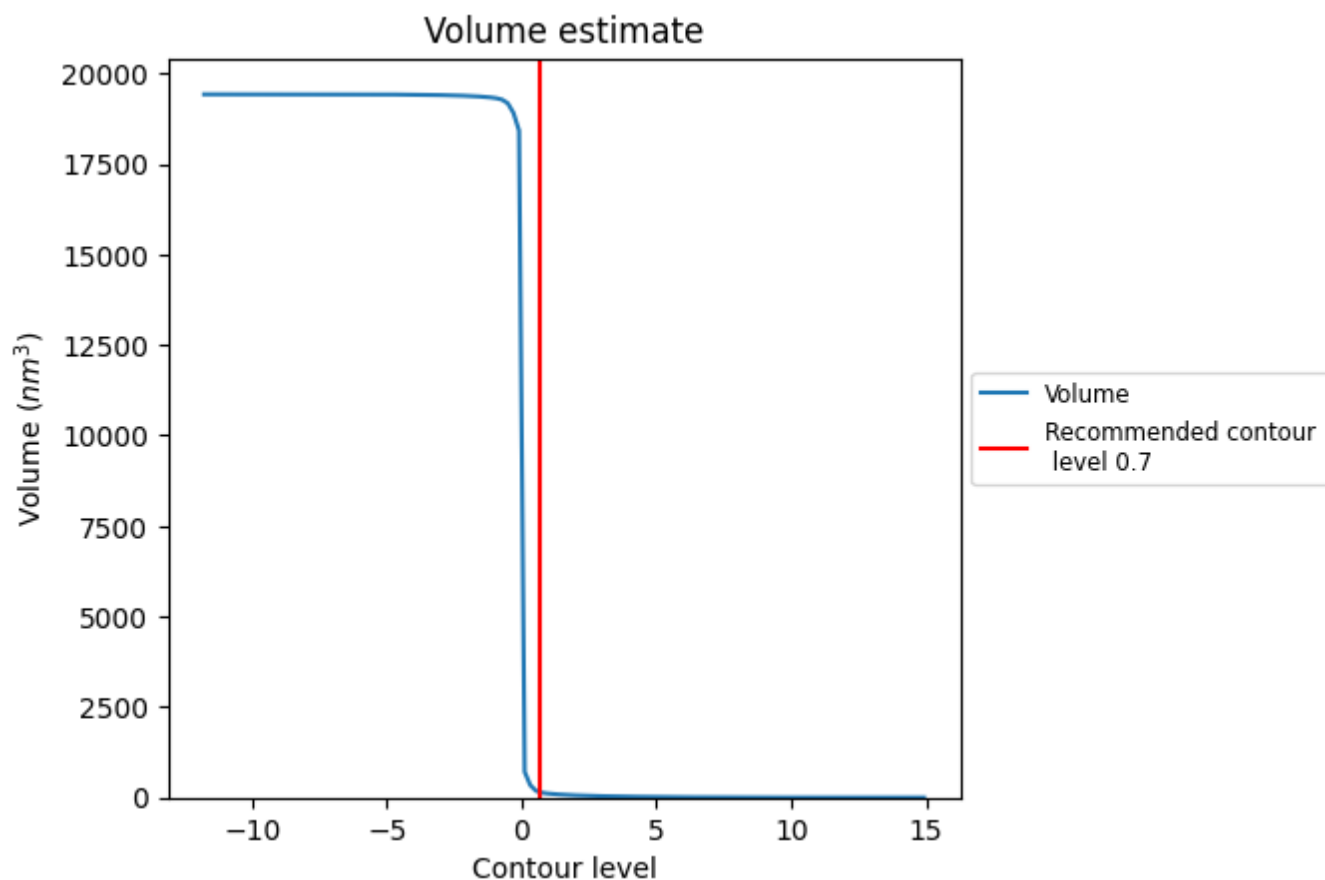
This section contains the results of statistical analysis of the map.

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.

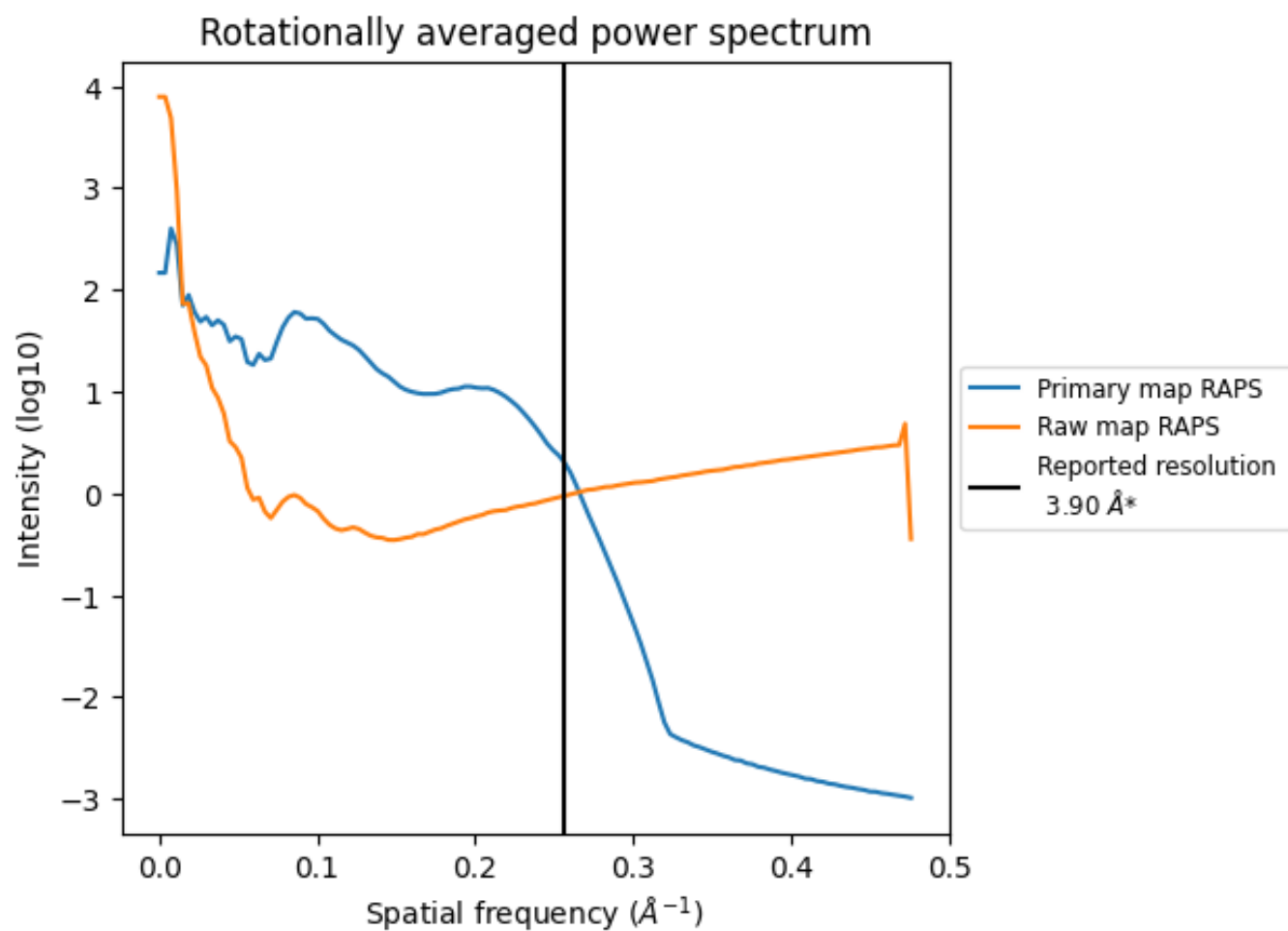
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 146 nm^3 ; this corresponds to an approximate mass of 132 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 ⓘ

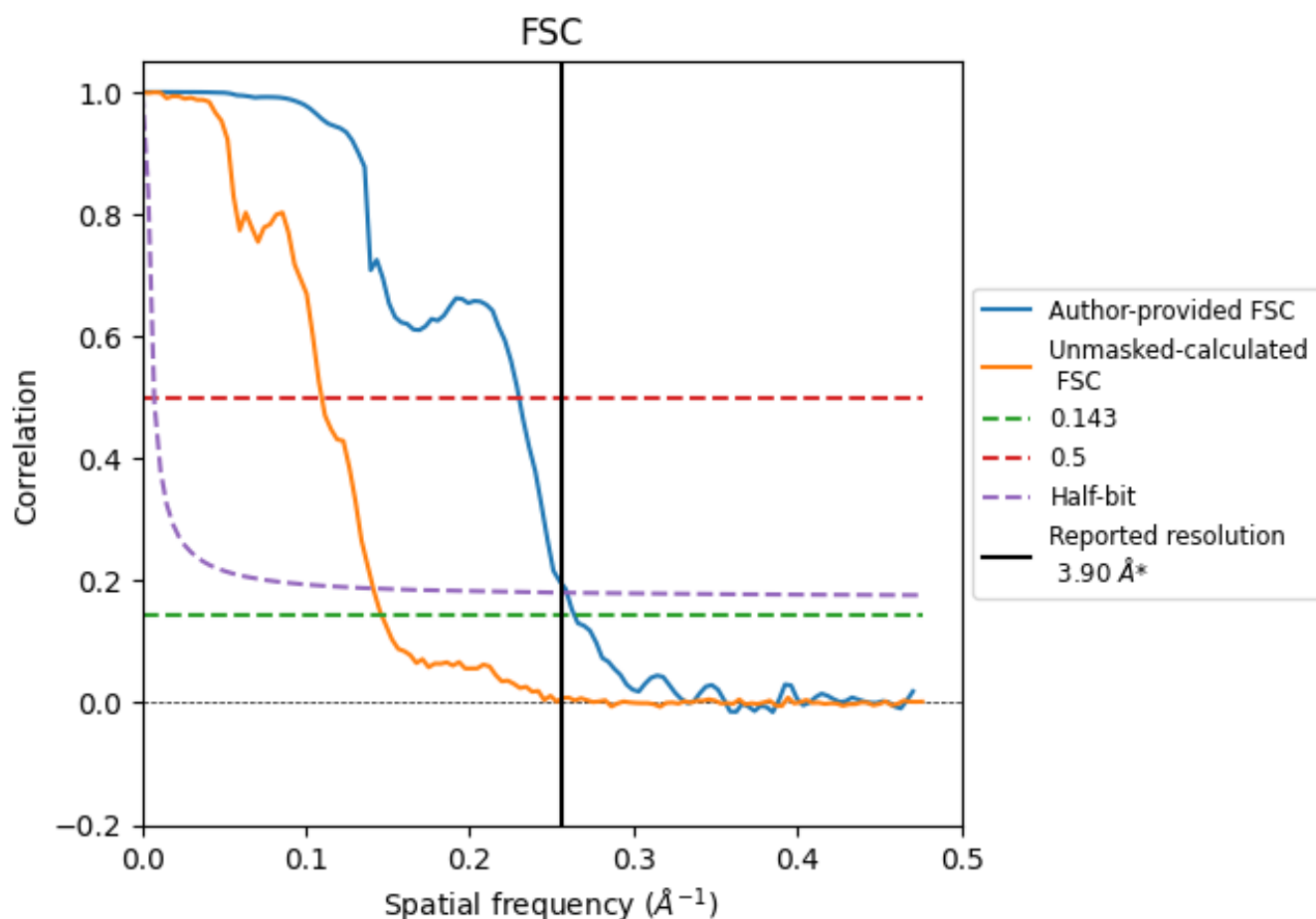


*Reported resolution corresponds to spatial frequency of $0.256 \text{ \AA}^{-1}$

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 0.256 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

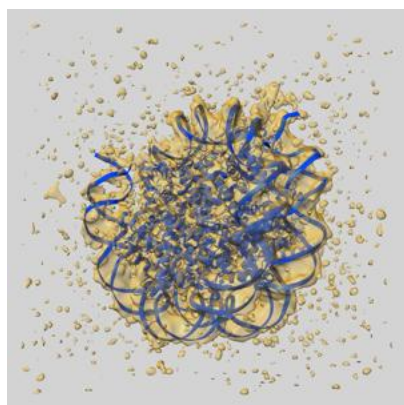
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.90	-	-
Author-provided FSC curve	3.79	4.35	3.86
Unmasked-calculated*	6.84	9.12	7.09

*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 6.84 differs from the reported value 3.9 by more than 10 %

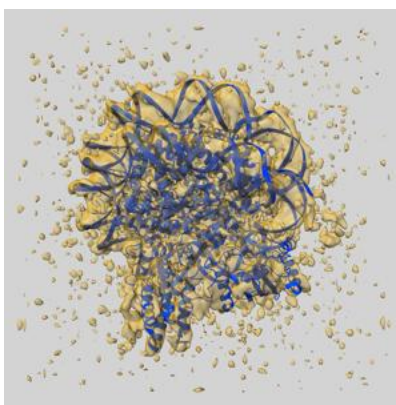
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-22581 and PDB model 7JZV. Per-residue inclusion information can be found in section [3](#) on page [7](#).

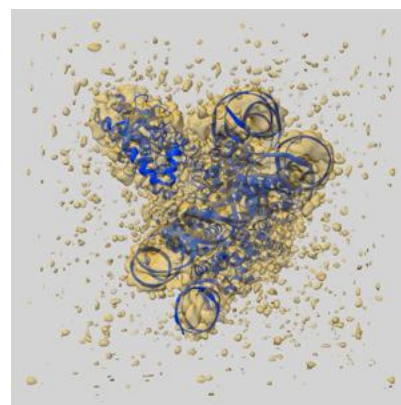
9.1 Map-model overlay [i](#)



X



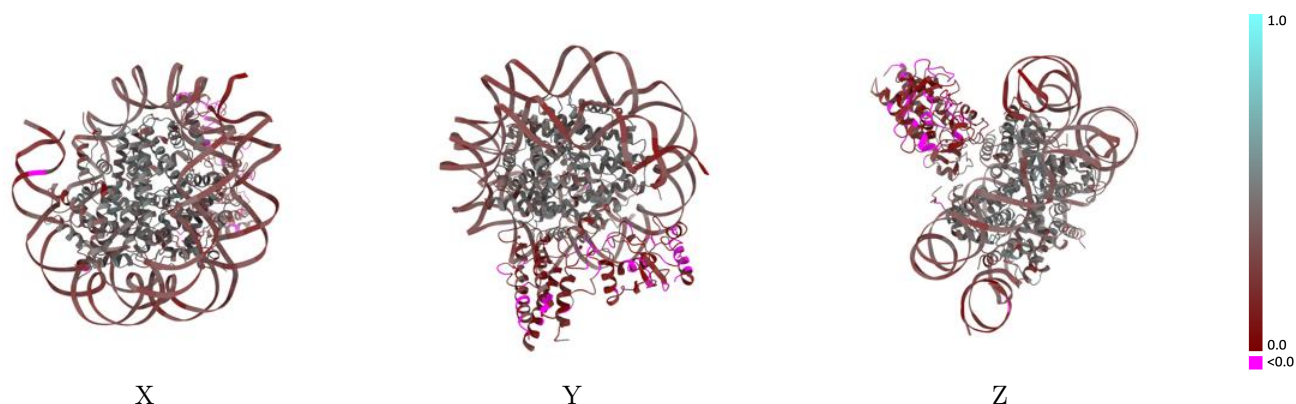
Y



Z

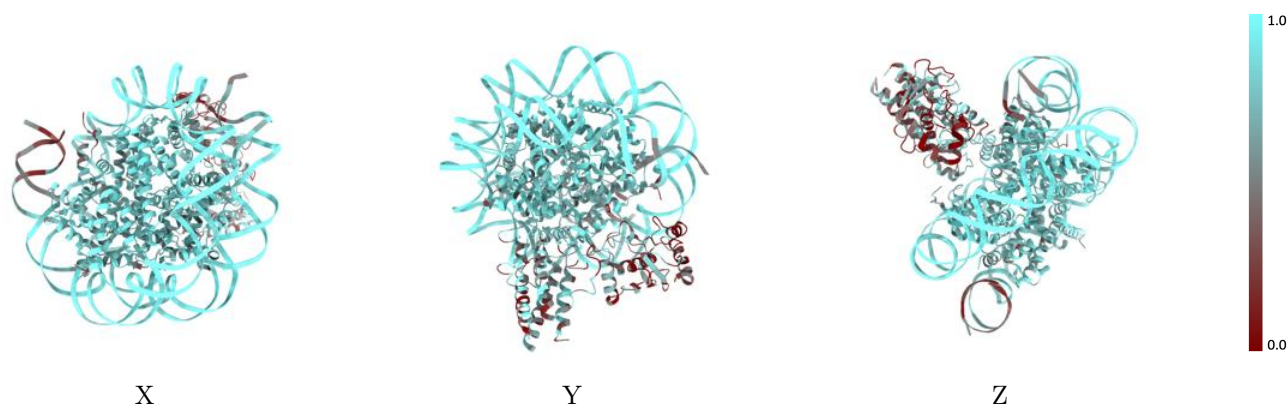
The images above show the 3D surface view of the map at the recommended contour level 0.7 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



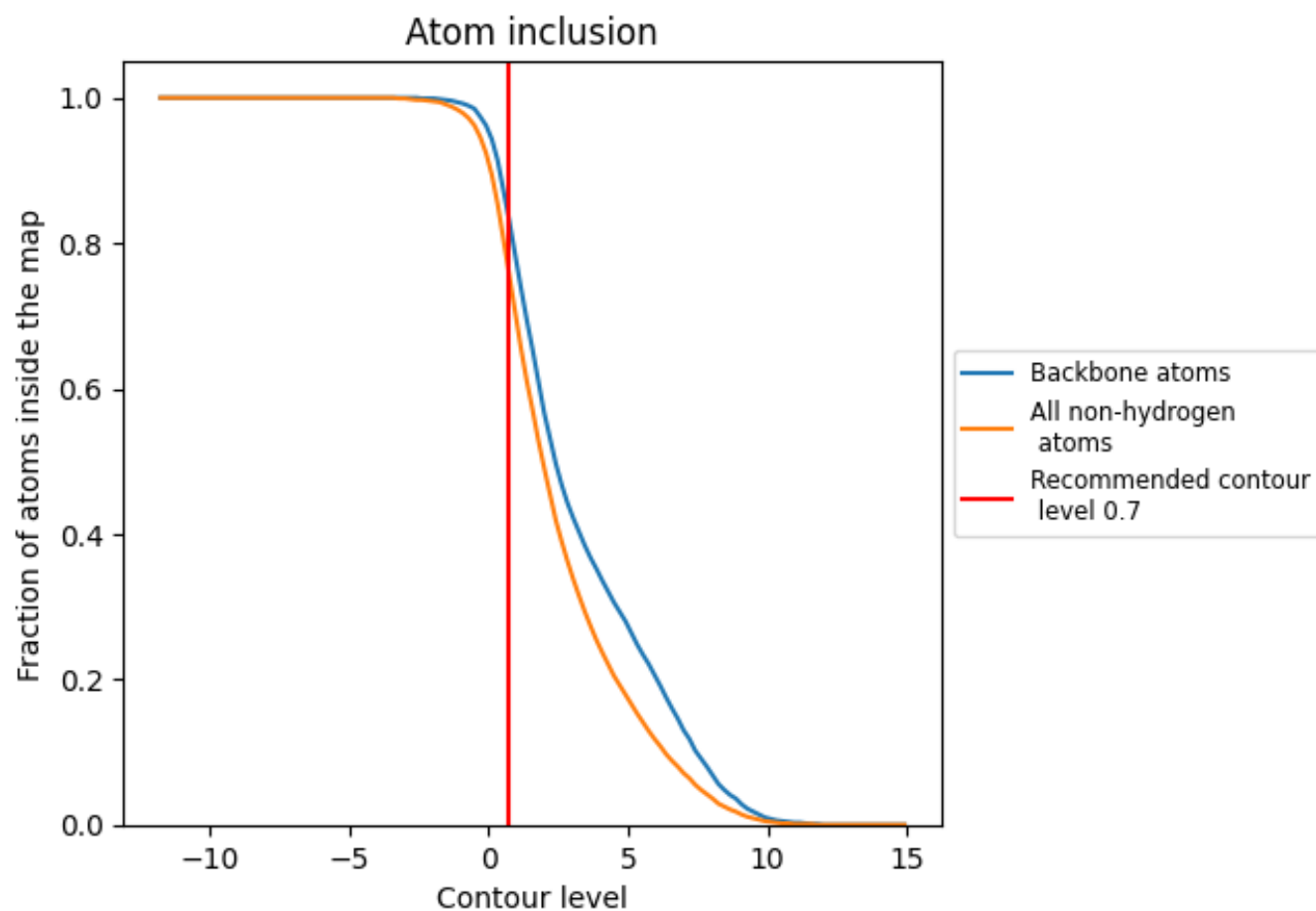
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.7).

9.4 Atom inclusion [i](#)



At the recommended contour level, 84% 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 (0.7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7660	0.3130
A	0.4580	0.1330
B	0.5650	0.1670
N	0.8180	0.4180
O	0.8130	0.4090
P	0.8270	0.4350
Q	0.8350	0.4240
X	0.8380	0.2890
Y	0.8460	0.2800
n	0.7720	0.3910
o	0.7930	0.3890
p	0.8530	0.4510
q	0.8840	0.4640

1.0

0.0

<0.0