



# wwPDB EM Validation Summary Report ⓘ

Jan 16, 2024 – 11:16 AM EST

PDB ID : 8TVQ  
EMDB ID : EMD-41648  
Title : Cryo-EM structure of CPD stalled 10-subunit Pol II in complex with Rad26  
Authors : Sarsam, R.D.; Lahiri, I.; Leschziner, A.E.  
Deposited on : 2023-08-18  
Resolution : 4.60 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](mailto: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>.

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

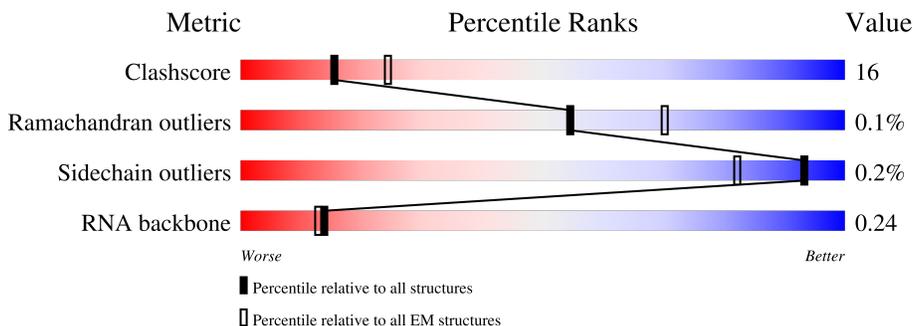
EMDB validation analysis : 0.0.1.dev70  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

# 1 Overall quality at a glance i

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

The reported resolution of this entry is 4.60 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	1733	
2	B	1224	
3	C	318	
4	E	215	
5	F	155	
6	H	146	
7	I	122	

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Mol	Chain	Length	Quality of chain
8	J	70	
9	K	120	
10	L	70	
11	M	434	
12	N	47	
13	T	46	
14	R	10	

## 2 Entry composition [i](#)

There are 16 unique types of molecules in this entry. The entry contains 57734 atoms, of which 26880 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 DNA-directed RNA polymerase II subunit RPB1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
1	A	1356	20724	6606	10250	1844	1965	59	0	0

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
2	B	1041	16000	5141	7876	1422	1508	53	0	0

- Molecule 3 is a protein called DNA-directed RNA polymerase II subunit RPB3.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
3	C	266	4148	1317	2053	348	417	13	0	0

- Molecule 4 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC1.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
4	E	215	3548	1116	1788	310	322	12	0	0

- Molecule 5 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC2.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
5	F	81	1198	394	578	110	114	2	0	0

- Molecule 6 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC3.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
6	H	133	2090	668	1030	174	213	5	0	0

- Molecule 7 is a protein called DNA-directed RNA polymerase II subunit RPB9.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
7	I	117	Total	C	H	N	O	S	0	0
			1138	451	396	142	144	5		

- Molecule 8 is a protein called DNA-directed RNA polymerases II subunit RPABC5.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
8	J	65	Total	C	H	N	O	S	0	0
			1074	339	542	93	94	6		

- Molecule 9 is a protein called DNA-directed RNA polymerase II subunit RPB11.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
9	K	104	Total	C	H	N	O	S	0	0
			1674	534	840	142	157	1		

- Molecule 10 is a protein called DNA-directed RNA polymerases II subunit RPABC4.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			S
10	L	46	Total	C	H	N	O	S	0	0
			751	224	387	72	64	4		

- Molecule 11 is a protein called DNA repair and recombination protein RAD26.

Mol	Chain	Residues	Atoms			AltConf	Trace	
			Total	C	N			O
11	M	434	Total	C	N	O	0	0
			2171	1302	434	435		

- Molecule 12 is a DNA chain called DNA (NTS).

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
12	N	46	Total	C	H	N	O	P	0	0
			1462	450	519	171	276	46		

- Molecule 13 is a DNA chain called DNA (TS).

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	H	N	O			P
13	T	45	Total	C	H	N	O	P	0	0
			1451	445	523	152	285	46		

- Molecule 14 is a RNA chain called RNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
			Total	C	H	N	O	P		
14	R	9	296	88	98	40	61	9	0	0

- Molecule 15 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
15	A	2	Total	Zn	0
			2	2	
15	B	1	Total	Zn	0
			1	1	
15	C	1	Total	Zn	0
			1	1	
15	I	2	Total	Zn	0
			2	2	
15	J	1	Total	Zn	0
			1	1	
15	L	1	Total	Zn	0
			1	1	

- Molecule 16 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

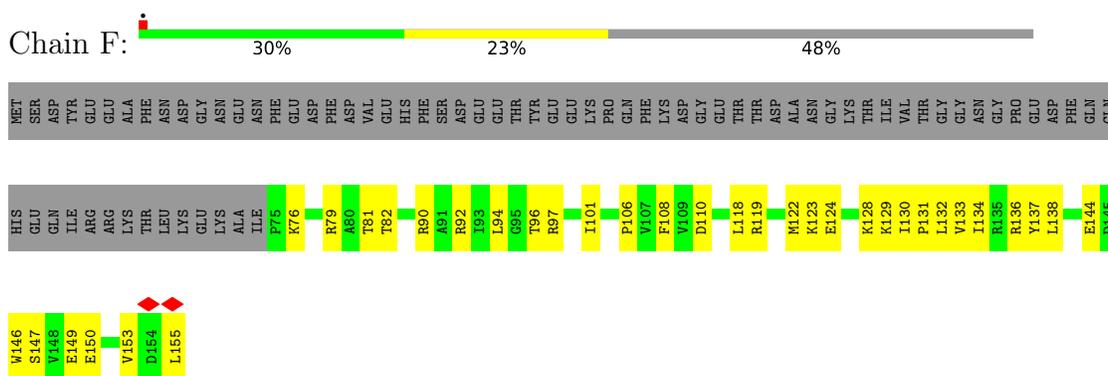
Mol	Chain	Residues	Atoms		AltConf
16	A	1	Total	Mg	0
			1	1	







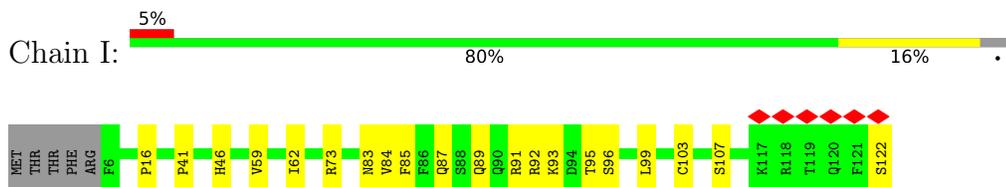
- Molecule 5: DNA-directed RNA polymerases I, II, and III subunit RPABC2



- Molecule 6: DNA-directed RNA polymerases I, II, and III subunit RPABC3



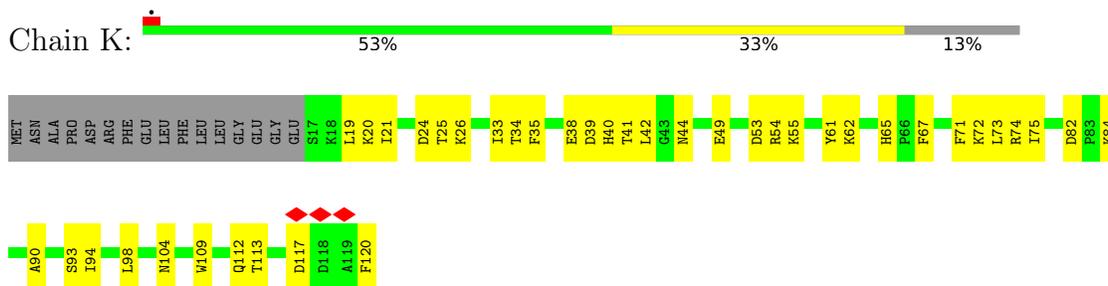
- Molecule 7: DNA-directed RNA polymerase II subunit RPB9



- Molecule 8: DNA-directed RNA polymerases II subunit RPABC5



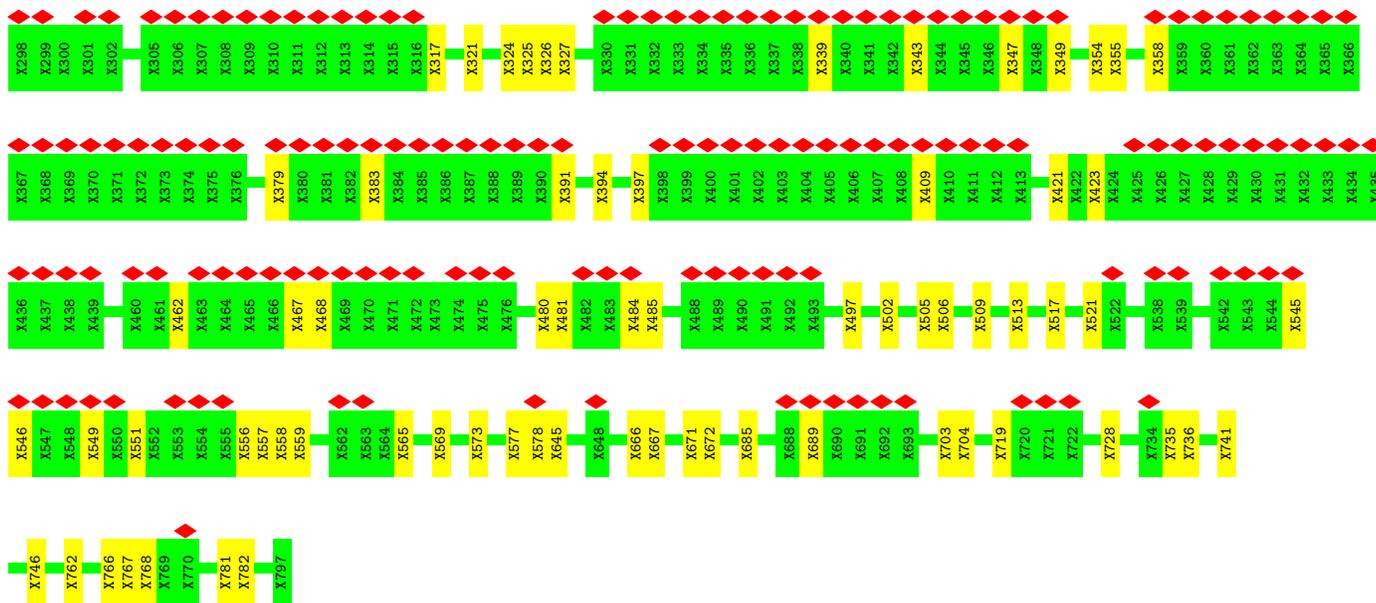
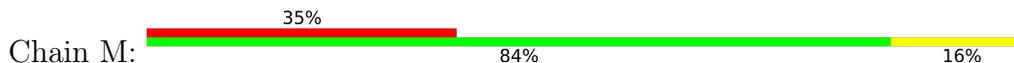
- Molecule 9: DNA-directed RNA polymerase II subunit RPB11



• Molecule 10: DNA-directed RNA polymerases II subunit RPABC4



• Molecule 11: DNA repair and recombination protein RAD26



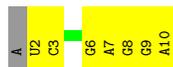
• Molecule 12: DNA (NTS)



• Molecule 13: DNA (TS)



• Molecule 14: RNA



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	22000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	52	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	2.691	Depositor
Minimum map value	-1.595	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.067	Depositor
Recommended contour level	0.26	Depositor
Map size (Å)	445.44, 445.44, 445.44	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.16, 1.16, 1.16	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: MG, ZN, TTD

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	0.28	0/10658	0.51	3/14420 (0.0%)
2	B	0.29	0/8277	0.50	1/11172 (0.0%)
3	C	0.29	0/2133	0.46	0/2891
4	E	0.28	0/1796	0.49	0/2416
5	F	0.29	0/630	0.60	1/851 (0.1%)
6	H	0.30	0/1076	0.51	0/1455
7	I	0.29	0/751	0.64	2/1024 (0.2%)
8	J	0.32	0/541	0.53	0/727
9	K	0.30	0/850	0.49	0/1148
10	L	0.28	0/366	0.58	0/485
12	N	0.38	0/1057	0.73	0/1629
13	T	0.66	0/989	0.96	2/1519 (0.1%)
14	R	0.76	0/222	0.85	0/345
All	All	0.32	0/29346	0.55	9/40082 (0.0%)

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	312	PRO	N-CA-CB	6.51	111.12	103.30
1	A	321	PRO	N-CA-CB	6.50	111.11	103.30
2	B	877	PRO	N-CA-CB	6.50	111.10	103.30
5	F	106	PRO	N-CA-CB	6.50	111.10	103.30
7	I	41	PRO	N-CA-CB	6.50	111.10	103.30

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	10474	10250	10393	359	0
2	B	8124	7876	7977	274	0
3	C	2095	2053	2053	74	0
4	E	1760	1788	1788	46	0
5	F	620	578	609	33	0
6	H	1060	1030	1030	42	0
7	I	742	396	534	15	0
8	J	532	542	545	33	0
9	K	834	840	839	38	0
10	L	364	387	388	12	0
11	M	2171	0	459	58	0
12	N	943	519	520	41	0
13	T	928	523	523	59	0
14	R	198	98	98	9	0
15	A	2	0	0	1	0
15	B	1	0	0	0	0
15	C	1	0	0	0	0
15	I	2	0	0	0	0
15	J	1	0	0	0	0
15	L	1	0	0	1	0
16	A	1	0	0	0	0
All	All	30854	26880	27756	961	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 16.

The worst 5 of 961 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)
11:M:484:UNK:CB	11:M:497:UNK:CB	1.74	1.56
11:M:355:UNK:CB	13:T:38:DA:OP2	1.81	1.26
11:M:317:UNK:CB	11:M:462:UNK:O	1.82	1.26
11:M:394:UNK:CB	13:T:38:DA:H5''	1.67	1.24
12:N:12:DC:N4	13:T:34:DT:N3	2.03	1.07

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	1342/1733 (77%)	1251 (93%)	89 (7%)	2 (0%)	51	85
2	B	1019/1224 (83%)	952 (93%)	67 (7%)	0	100	100
3	C	264/318 (83%)	254 (96%)	10 (4%)	0	100	100
4	E	213/215 (99%)	202 (95%)	11 (5%)	0	100	100
5	F	79/155 (51%)	74 (94%)	5 (6%)	0	100	100
6	H	127/146 (87%)	116 (91%)	11 (9%)	0	100	100
7	I	115/122 (94%)	106 (92%)	9 (8%)	0	100	100
8	J	63/70 (90%)	58 (92%)	5 (8%)	0	100	100
9	K	102/120 (85%)	99 (97%)	3 (3%)	0	100	100
10	L	44/70 (63%)	40 (91%)	4 (9%)	0	100	100
All	All	3368/4173 (81%)	3152 (94%)	214 (6%)	2 (0%)	54	85

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1098	VAL
1	A	958	VAL

### 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	1124/1520 (74%)	1122 (100%)	2 (0%)	93	96
2	B	856/1061 (81%)	853 (100%)	3 (0%)	91	94
3	C	234/274 (85%)	234 (100%)	0	100	100
4	E	197/197 (100%)	196 (100%)	1 (0%)	88	93
5	F	60/137 (44%)	60 (100%)	0	100	100
6	H	117/128 (91%)	117 (100%)	0	100	100
7	I	50/116 (43%)	50 (100%)	0	100	100
8	J	60/65 (92%)	60 (100%)	0	100	100
9	K	89/102 (87%)	89 (100%)	0	100	100
10	L	40/57 (70%)	40 (100%)	0	100	100
All	All	2827/3657 (77%)	2821 (100%)	6 (0%)	93	96

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	605	ARG
2	B	995	ARG
4	E	192	ARG
1	A	1194	ARG
1	A	840	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	800	GLN
2	B	1076	HIS
3	C	65	HIS
2	B	1112	GLN
1	A	510	GLN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
14	R	8/10 (80%)	1 (12%)	1 (12%)

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
14	R	3	C

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
14	R	8	G

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

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
13	TTD	T	18	13	42,45,46	3.30	16 (38%)	62,74,77	2.14	22 (35%)

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
13	TTD	T	18	13	-	12/22/109/110	0/5/6/6

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	T	18	TTD	C5-C6	-8.45	1.45	1.55
13	T	18	TTD	C2'-C3R	-7.47	1.36	1.52
13	T	18	TTD	C5T-C6T	-7.40	1.46	1.55
13	T	18	TTD	C2-N3	6.71	1.49	1.38
13	T	18	TTD	C2T-N3T	6.08	1.48	1.38

The worst 5 of 22 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	T	18	TTD	C2'-C1'-N1	-5.44	108.24	115.59
13	T	18	TTD	C4T-N3T-C2T	-4.71	119.44	126.67
13	T	18	TTD	N3T-C2T-N1T	4.66	121.53	116.69
13	T	18	TTD	C4-N3-C2	-4.57	119.65	126.67
13	T	18	TTD	O4T-C4T-C5T	-3.92	119.75	122.88

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
13	T	18	TTD	O4'-C1'-N1-C2
13	T	18	TTD	C5R-O5R-PB-O5P
13	T	18	TTD	C5R-O5R-PB-O4P
13	T	18	TTD	C3'-C4'-C5R-O5R
13	T	18	TTD	O4'-C4R-C5'-O5'

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	T	18	TTD	7	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 9 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
11	M	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	M	388:UNK	C	389:UNK	N	3.71
1	M	578:UNK	C	645:UNK	N	3.51

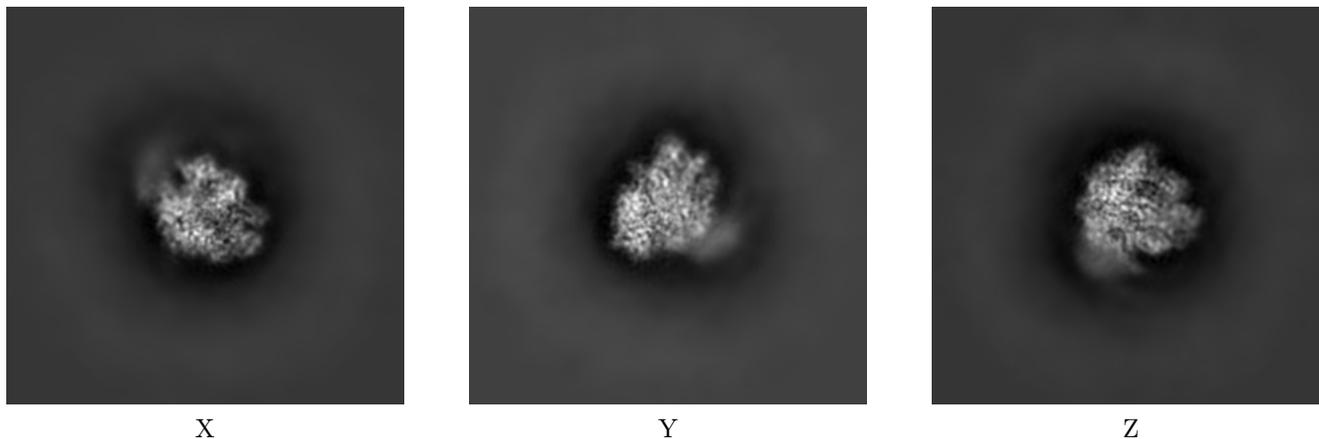
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-41648. These allow visual inspection of the internal detail of the map and identification of artifacts.

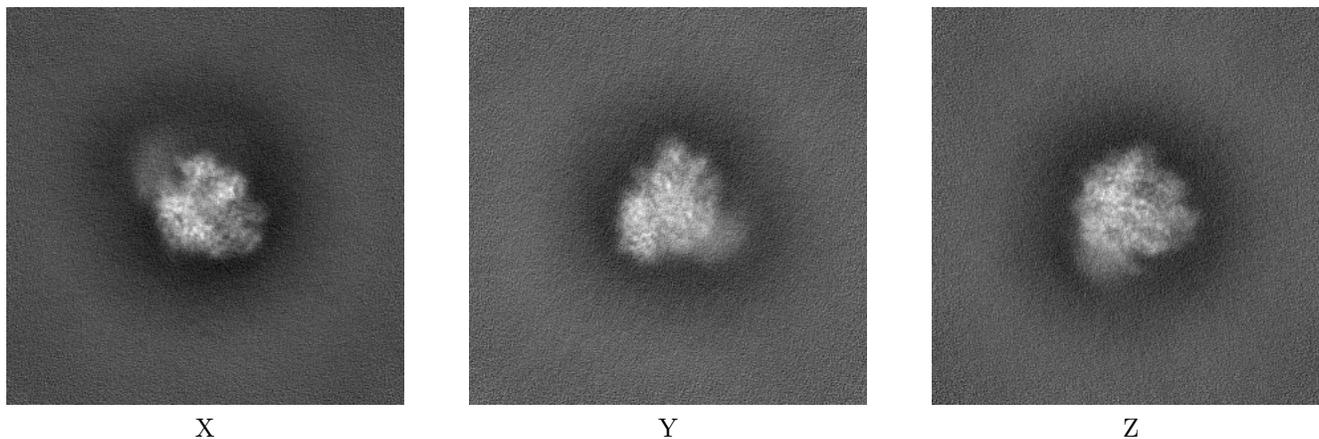
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

#### 6.1.1 Primary map



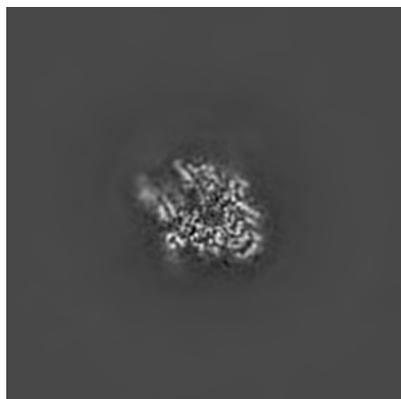
#### 6.1.2 Raw map



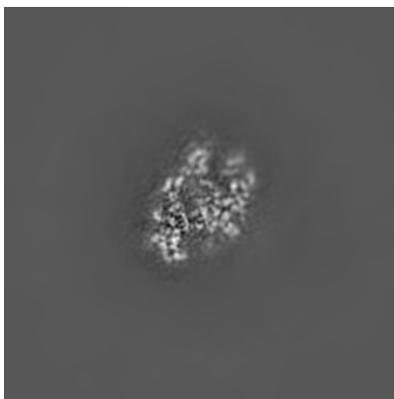
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

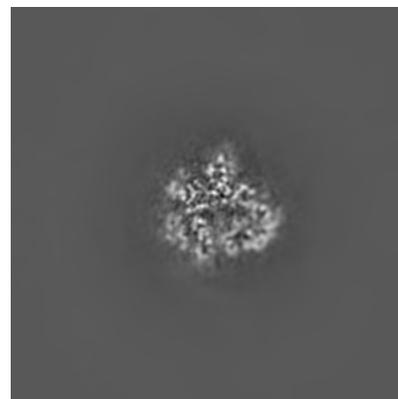
### 6.2.1 Primary map



X Index: 192

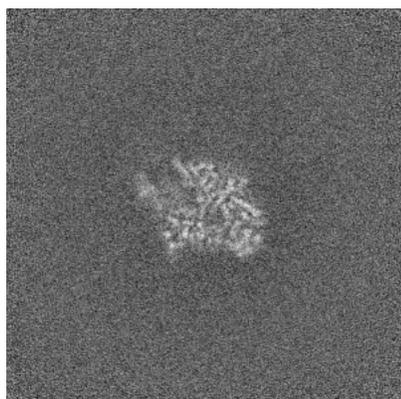


Y Index: 192

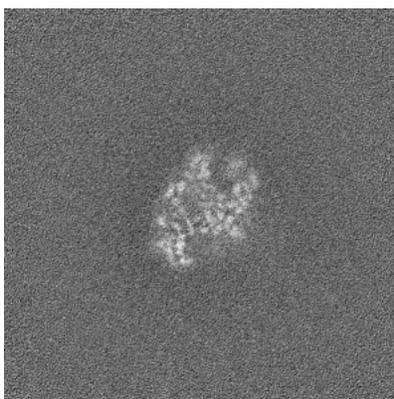


Z Index: 192

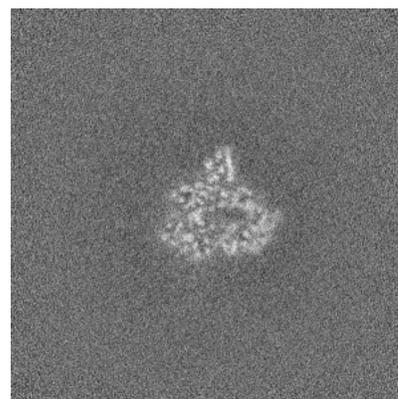
### 6.2.2 Raw map



X Index: 192



Y Index: 192

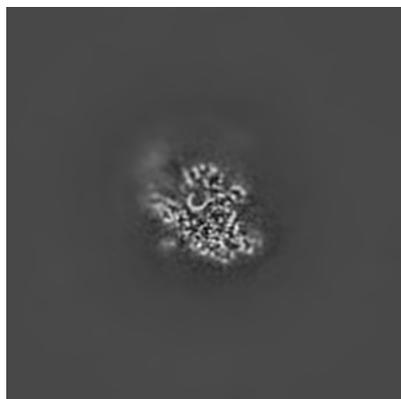


Z Index: 192

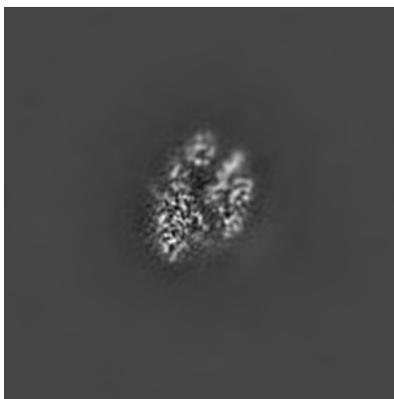
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

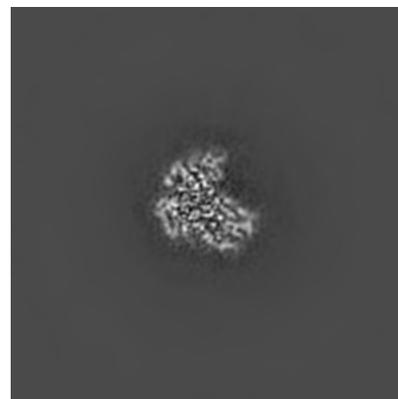
### 6.3.1 Primary map



X Index: 181

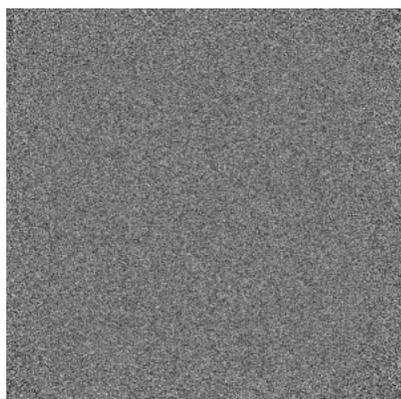


Y Index: 185

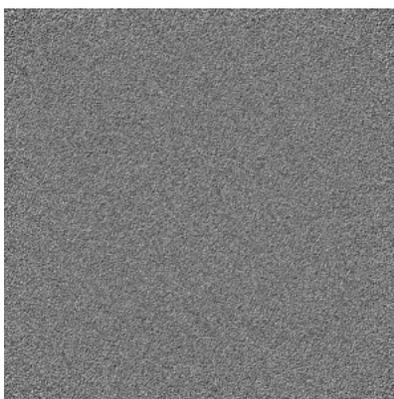


Z Index: 164

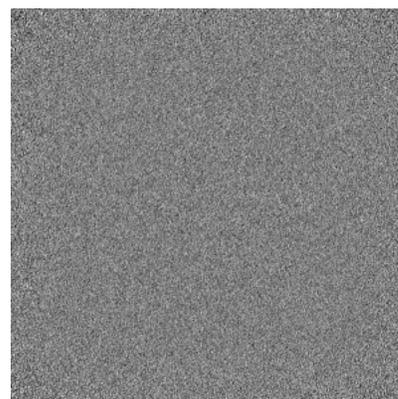
### 6.3.2 Raw map



X Index: 0



Y Index: 0

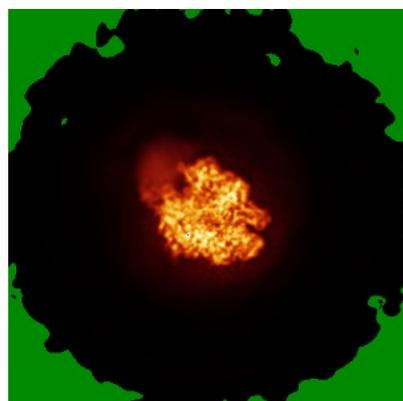


Z Index: 0

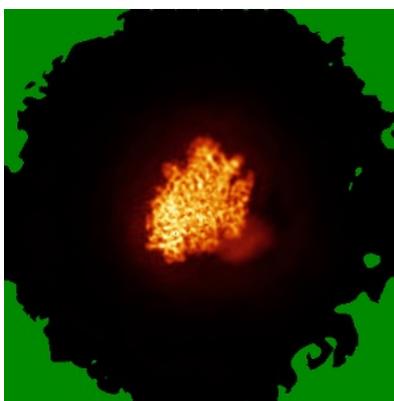
The images above show the largest variance slices of the map in three orthogonal directions.

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

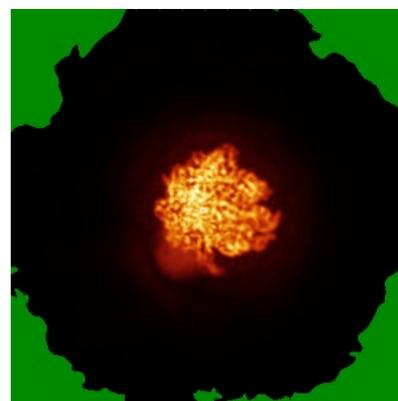
### 6.4.1 Primary map



X

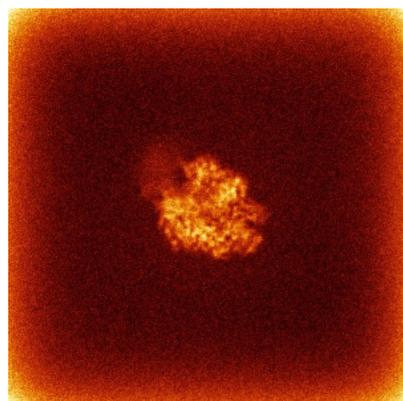


Y

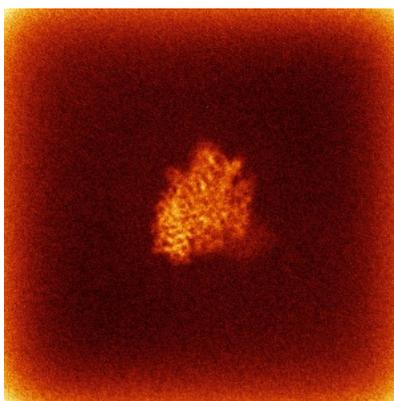


Z

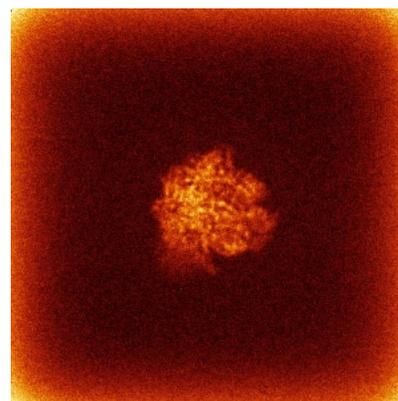
### 6.4.2 Raw map



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.

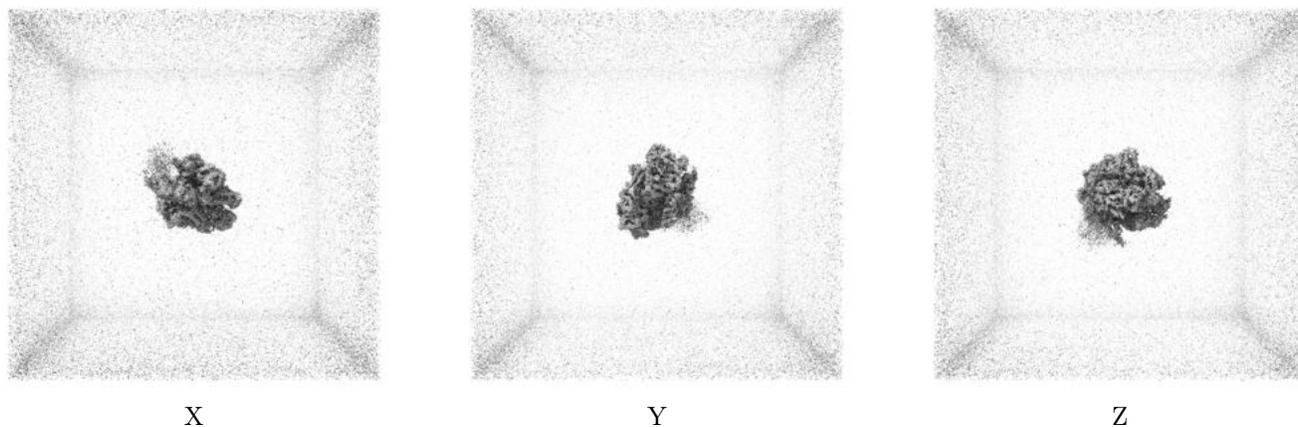
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.26. 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.

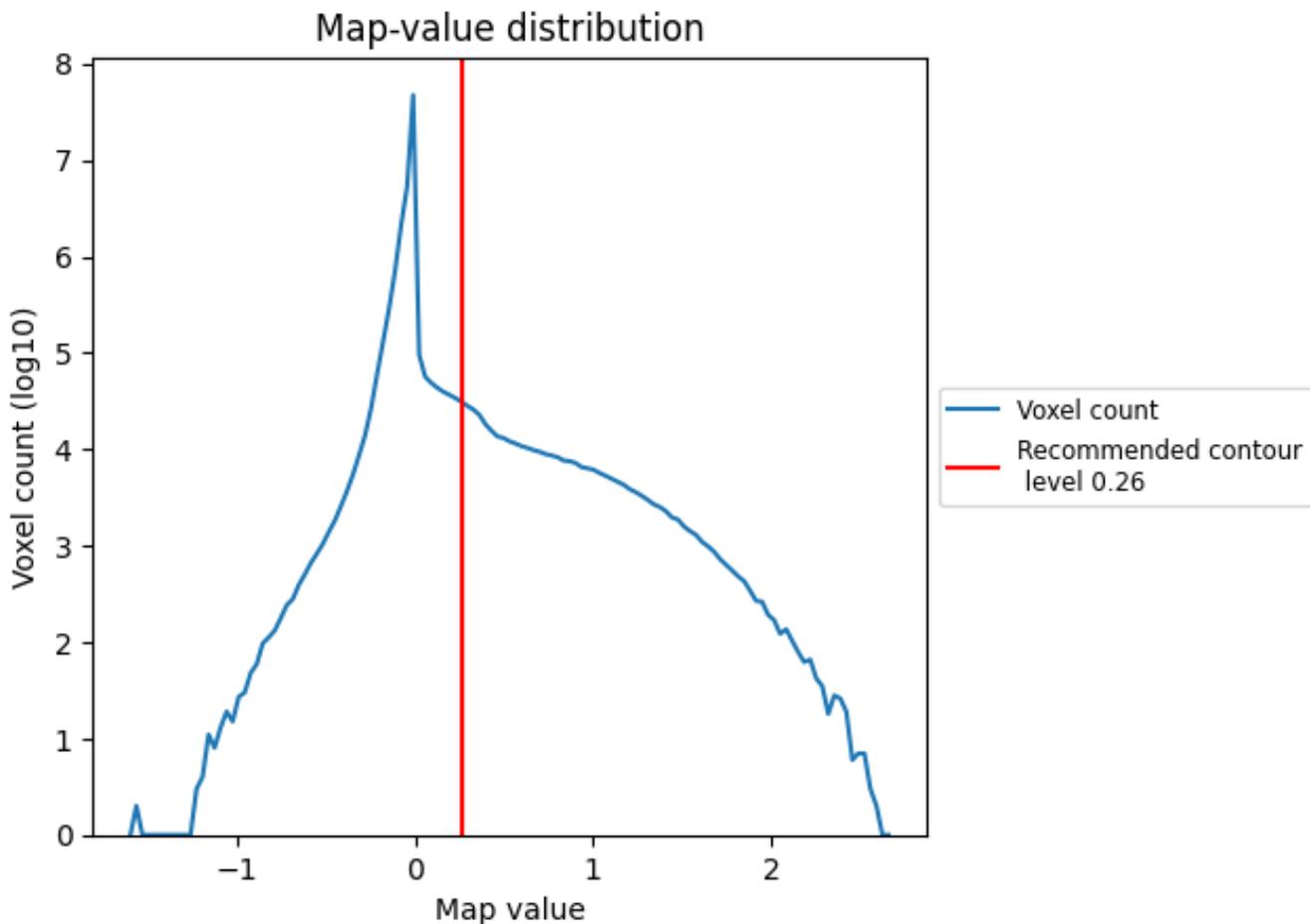
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

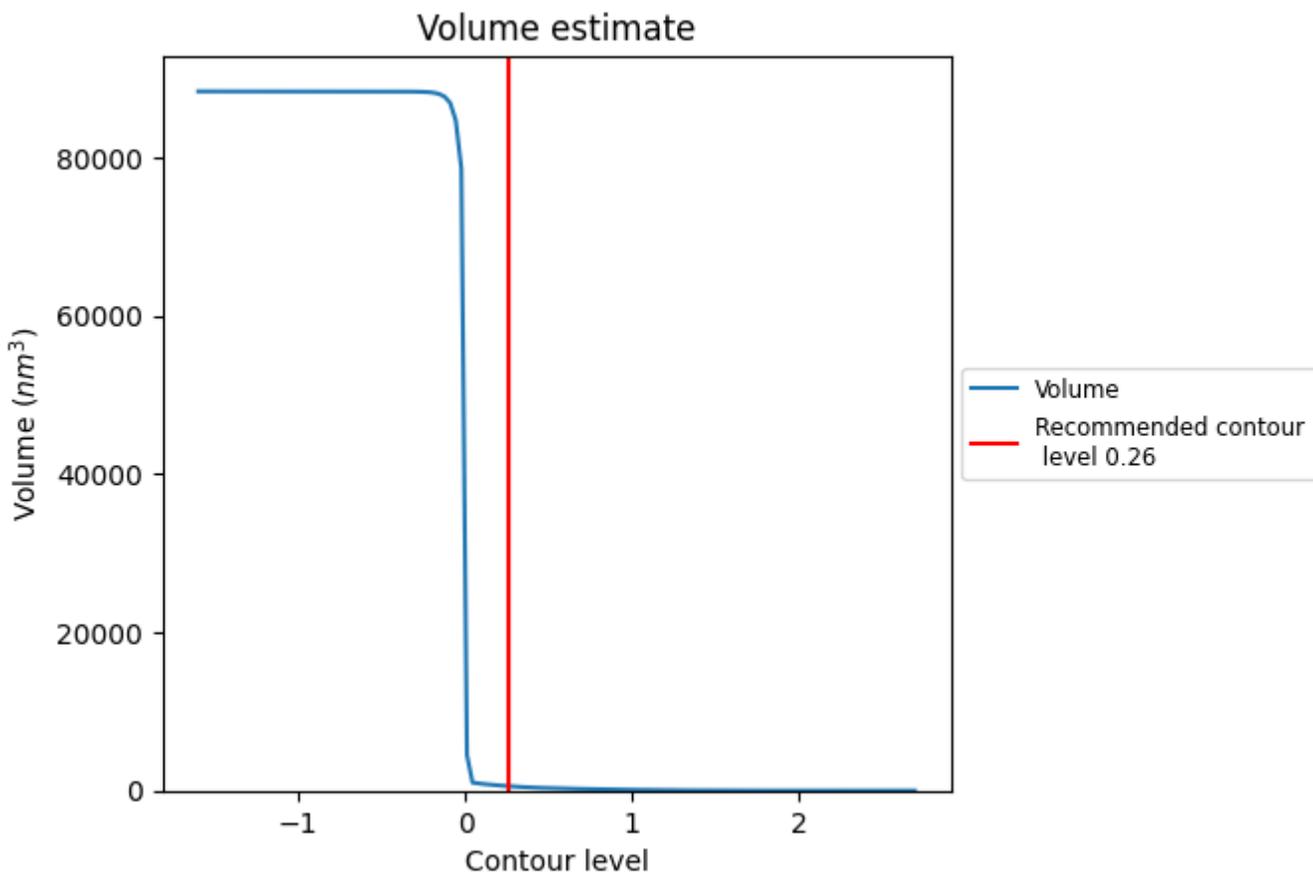
This section contains the results of statistical analysis of the map.

### 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.

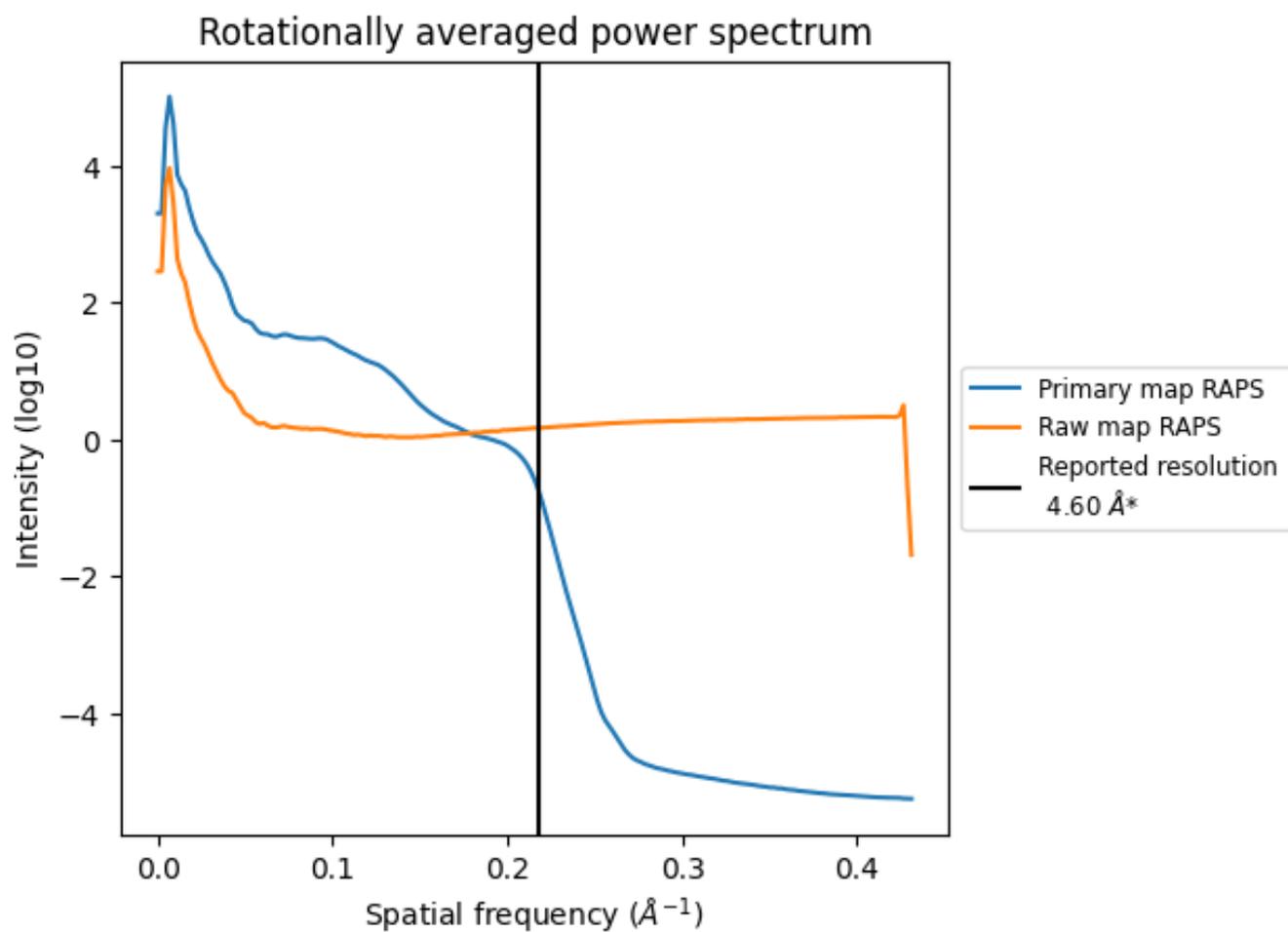
## 7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 566 nm<sup>3</sup>; this corresponds to an approximate mass of 511 kDa.

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

### 7.3 Rotationally averaged power spectrum i

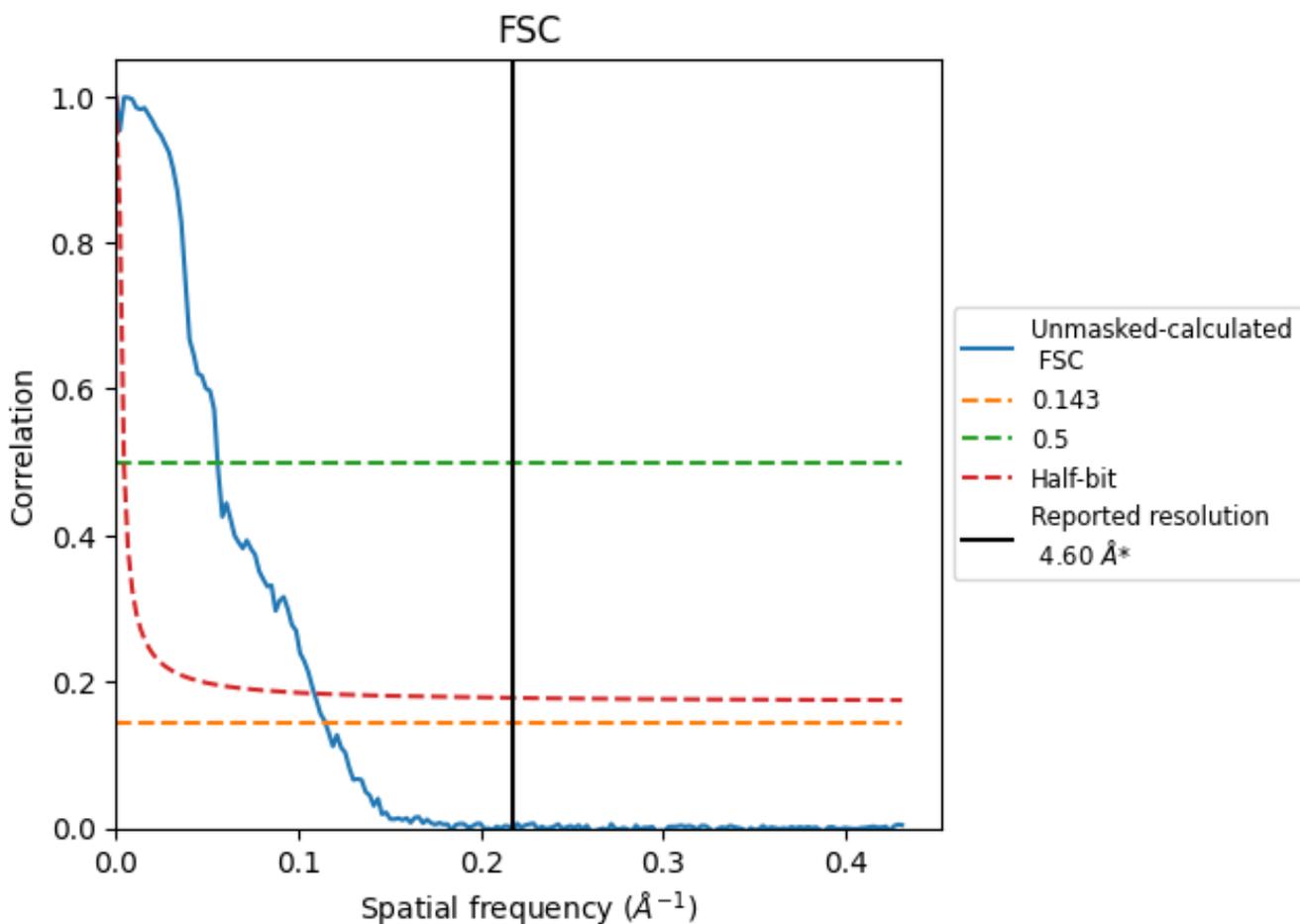


\*Reported resolution corresponds to spatial frequency of 0.217 Å<sup>-1</sup>

## 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 0.217  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

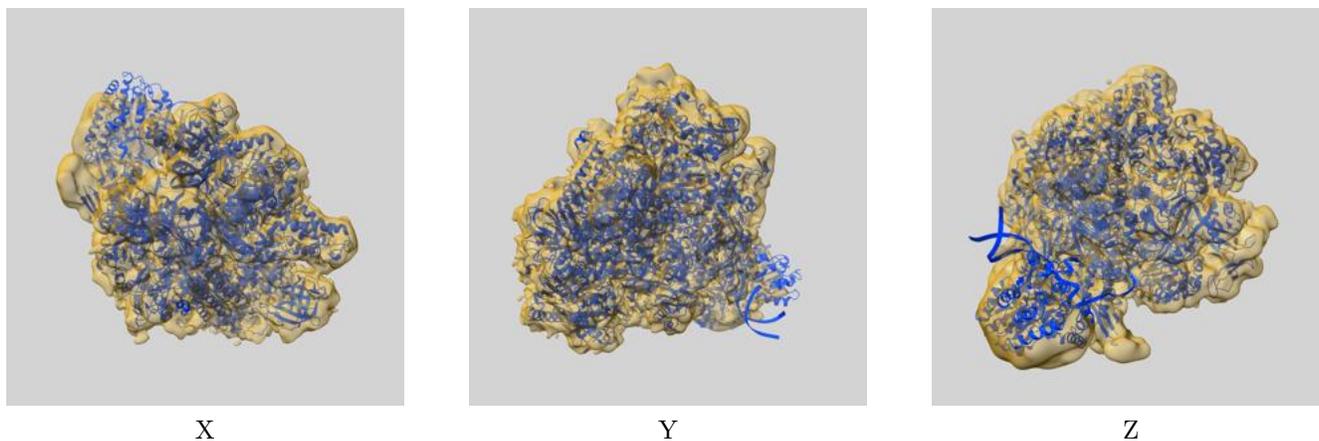
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.60	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	8.70	17.92	9.19

\*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 8.70 differs from the reported value 4.6 by more than 10 %

## 9 Map-model fit [i](#)

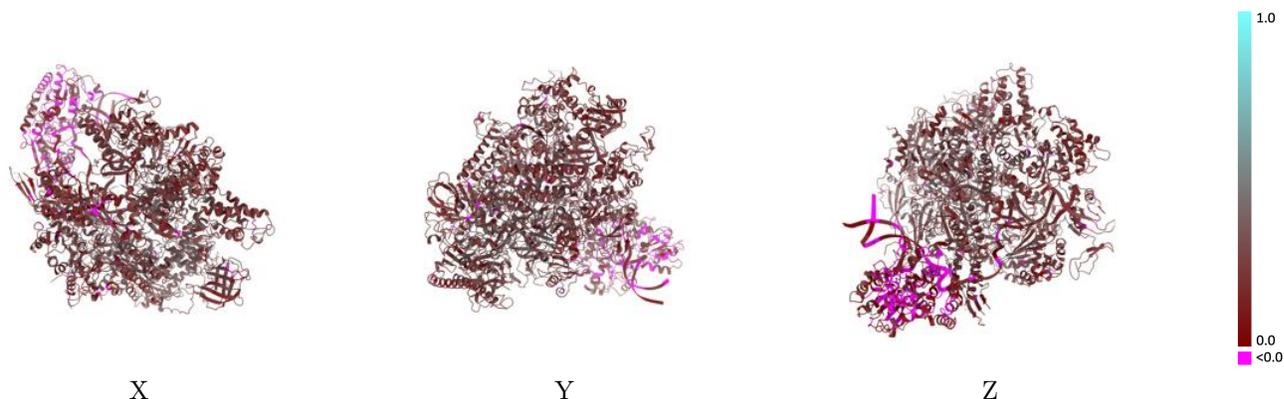
This section contains information regarding the fit between EMDB map EMD-41648 and PDB model 8TVQ. Per-residue inclusion information can be found in section 3 on page 7.

### 9.1 Map-model overlay [i](#)



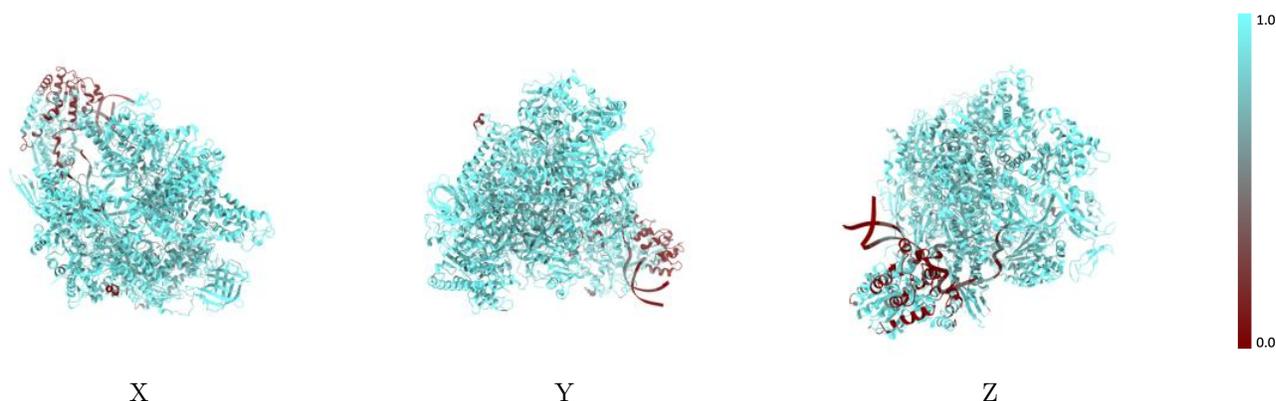
The images above show the 3D surface view of the map at the recommended contour level 0.26 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



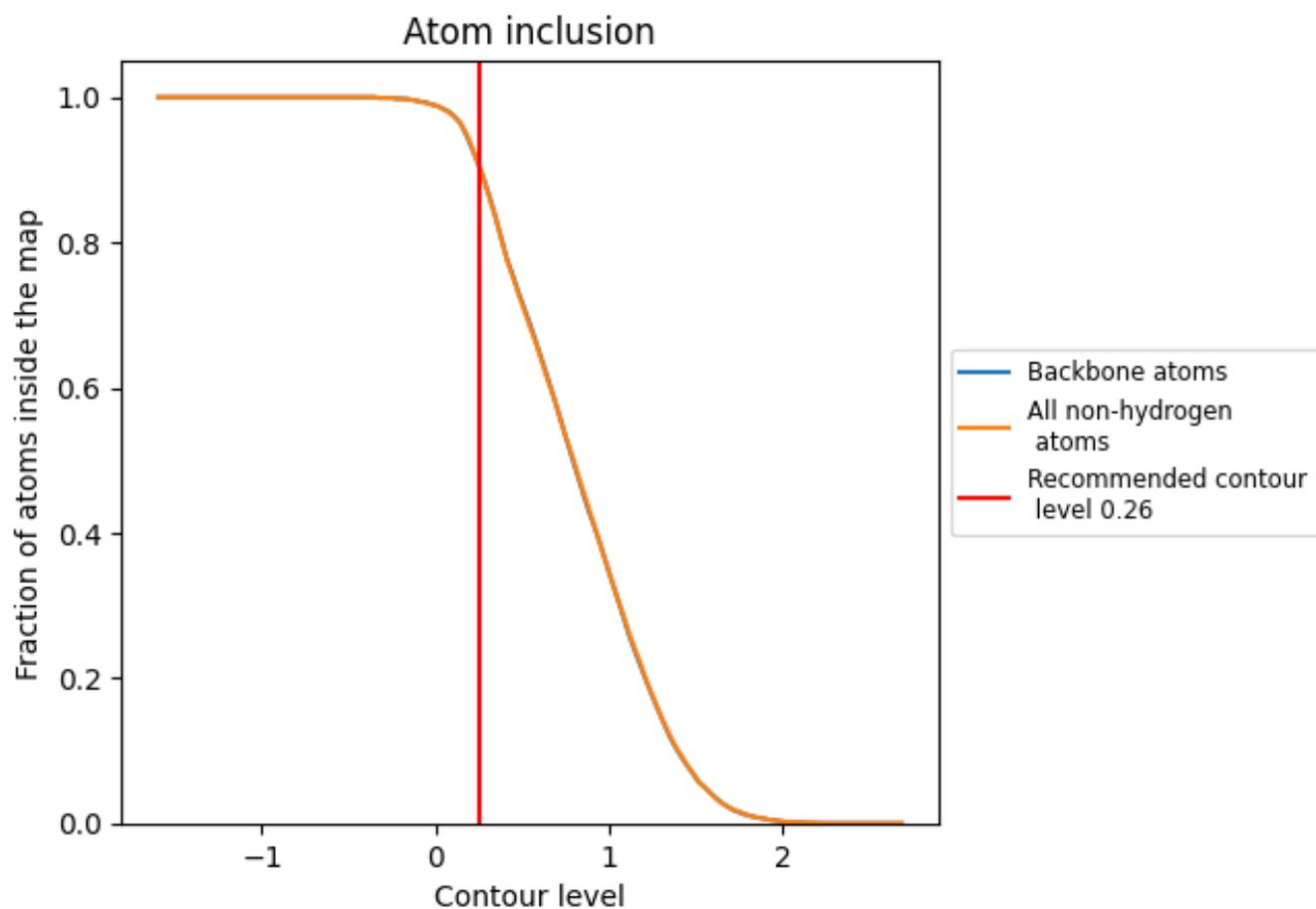
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.26).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 90% of all backbone atoms, 90% 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 (0.26) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9010	 0.2470
A	 0.9340	 0.2690
B	 0.9290	 0.2730
C	 0.9640	 0.2900
E	 0.9740	 0.2430
F	 0.9290	 0.2590
H	 0.9840	 0.2450
I	 0.9470	 0.2500
J	 0.9790	 0.3050
K	 0.9500	 0.2690
L	 0.9690	 0.2540
M	 0.6560	 0.0610
N	 0.5240	 0.0990
R	 0.9800	 0.2890
T	 0.7520	 0.1800

