



Full wwPDB EM Validation Report ⓘ

Jun 9, 2026 – 02:28 pm BST

PDB ID : 9RYB / pdb_00009ryb
EMDB ID : EMD-54374
Title : CryoEM structure of transcribing RNA polymerase II elongation complex in post-catalysis state
Authors : Li, Q.; Yi, G.; Zhang, P.; Wang, D.
Deposited on : 2025-07-15
Resolution : 2.33 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:

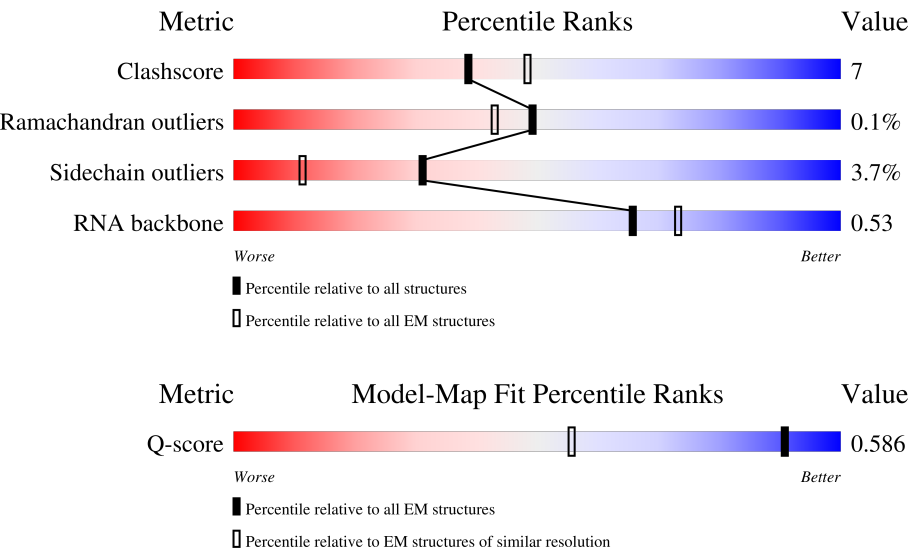
EMDB validation analysis : 0.0.1.dev132
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.33 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	4434 (1.83 - 2.83)

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	1733	8% 64% 16% • 19%
2	B	1224	10% 78% 16% • 5%
3	C	318	• 70% 13% • 16%

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Mol	Chain	Length	Quality of chain
4	D	221	76% 61% 14% 24%
5	E	215	14% 88% 11%
6	F	155	39% 14% 46%
7	G	171	92% 74% 25%
8	H	146	81% 11% 8%
9	I	122	55% 62% 28% 7%
10	J	70	79% 13% 7%
11	K	120	87% 6% 7%
12	L	70	13% 49% 16% 36%
13	N	74	16% 7% 15% 78%
14	R	19	26% 16% 16% 42%
15	T	74	14% 16% 20% 64%

2 Entry composition

There are 20 unique types of molecules in this entry. The entry contains 33089 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 DNA-directed RNA polymerase II subunit RPB1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1397	Total	C	N	O	S	0	0
			10999	6941	1916	2080	62		

- Molecule 2 is a protein called DNA-directed RNA polymerase II subunit RPB2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	1164	Total	C	N	O	S	0	0
			9268	5855	1627	1731	55		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	266	Total	C	N	O	S	0	0
			2095	1317	348	417	13		

- Molecule 4 is a protein called DNA-directed RNA polymerase II subunit RPB4.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	167	Total	C	N	O	S	0	0
			1327	821	230	273	3		

- Molecule 5 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	213	Total	C	N	O	S	0	0
			1744	1107	308	318	11		

- Molecule 6 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	83	Total	C	N	O	S	0	0
			670	428	114	125	3		

- Molecule 7 is a protein called DNA-directed RNA polymerase II subunit RPB7.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	171	Total	C	N	O	S	0	0
			1340	861	222	249	8		

- Molecule 8 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC3.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	134	Total	C	N	O	S	0	0
			1077	679	182	212	4		

- Molecule 9 is a protein called DNA-directed RNA polymerase II subunit RPB9.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	114	Total	C	N	O	S	0	0
			926	570	167	178	11		

- Molecule 10 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC5.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	65	Total	C	N	O	S	0	0
			532	339	93	94	6		

- Molecule 11 is a protein called DNA-directed RNA polymerase II subunit RPB11.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	112	Total	C	N	O	S	0	0
			904	580	154	168	2		

- Molecule 12 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC4.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	45	Total	C	N	O	S	0	0
			359	221	71	63	4		

- Molecule 13 is a DNA chain called DNA (74-mer).

Mol	Chain	Residues	Atoms					AltConf	Trace
13	N	16	Total	C	N	O	P	0	0
			334	156	69	93	16		

- Molecule 14 is a RNA chain called RNA (5'-R(*AP*UP*CP*GP*AP*GP*AP*GP*GP*AP

*UP*GP*GP*GP*AP*GP*AP*AP*G)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
14	R	11	Total	C	N	O	P	0	0
			246	109	52	74	11		

- Molecule 15 is a DNA chain called DNA (74-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
15	T	27	Total	C	N	O	P	0	0
			537	257	85	168	27		

- Molecule 16 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

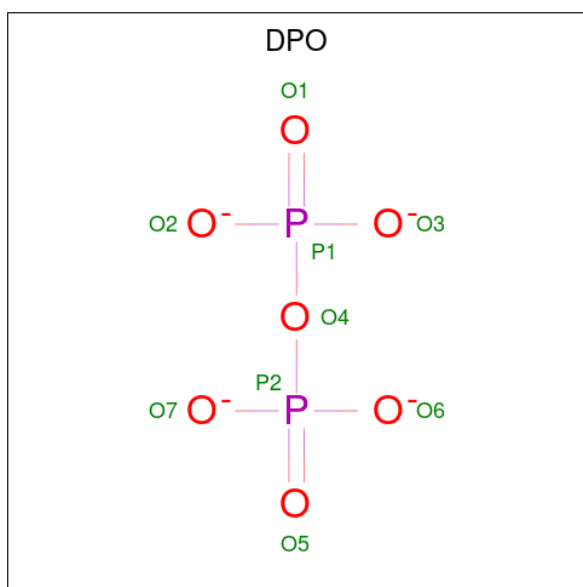
- Molecule 17 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 18 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
18	A	1	Total	K	0
			1	1	

- Molecule 19 is DIPHOSPHATE (CCD ID: DPO) (formula: O₇P₂).



Mol	Chain	Residues	Atoms			AltConf
19	B	1	Total	O	P	0
			9	7	2	

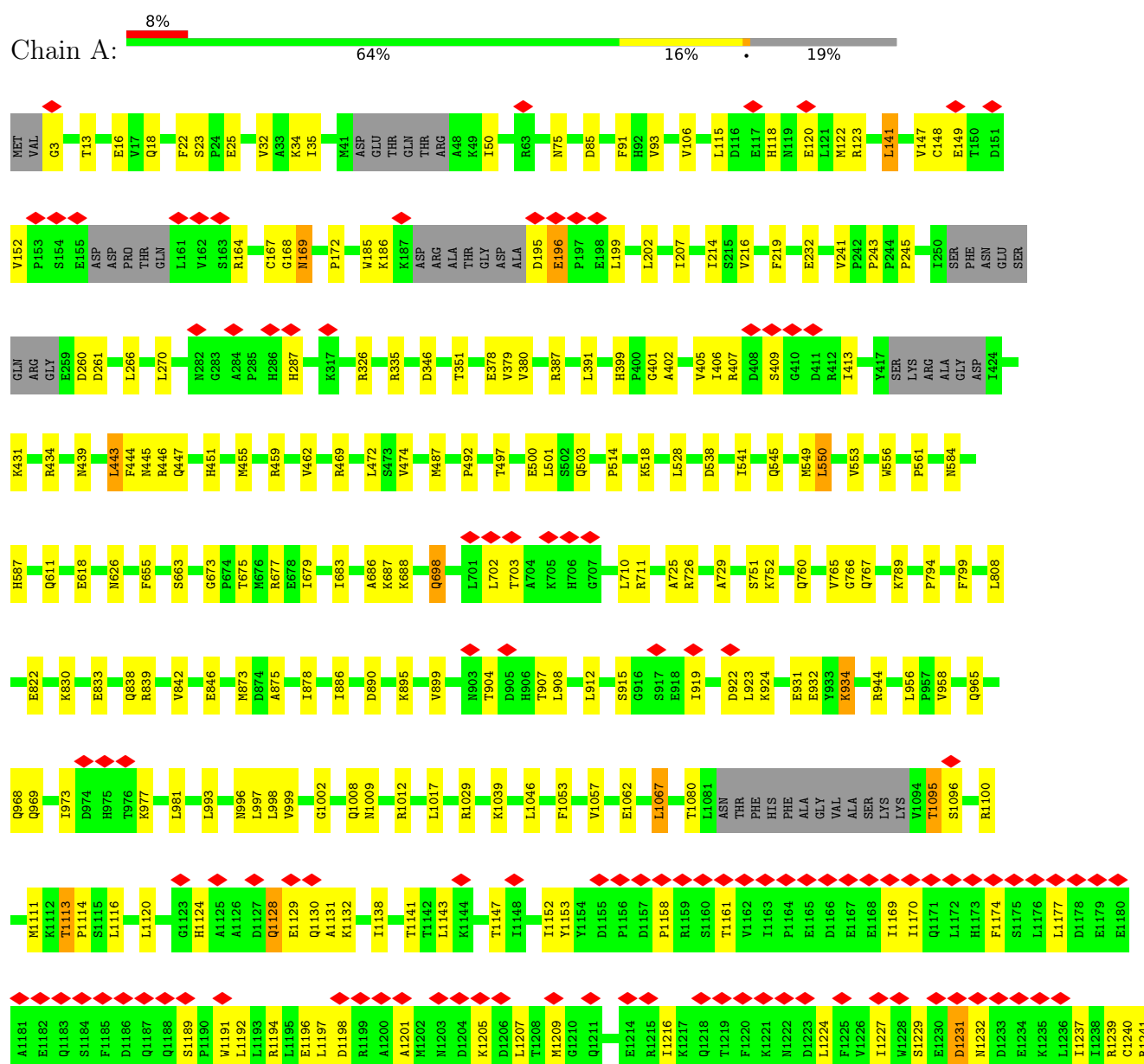
- Molecule 20 is water.

Mol	Chain	Residues	Atoms		AltConf
20	A	267	Total	O	0
			267	267	
20	B	232	Total	O	0
			232	232	
20	C	68	Total	O	0
			68	68	
20	E	22	Total	O	0
			22	22	
20	F	15	Total	O	0
			15	15	
20	H	10	Total	O	0
			10	10	
20	J	18	Total	O	0
			18	18	
20	K	30	Total	O	0
			30	30	
20	L	4	Total	O	0
			4	4	
20	R	22	Total	O	0
			22	22	
20	T	24	Total	O	0
			24	24	

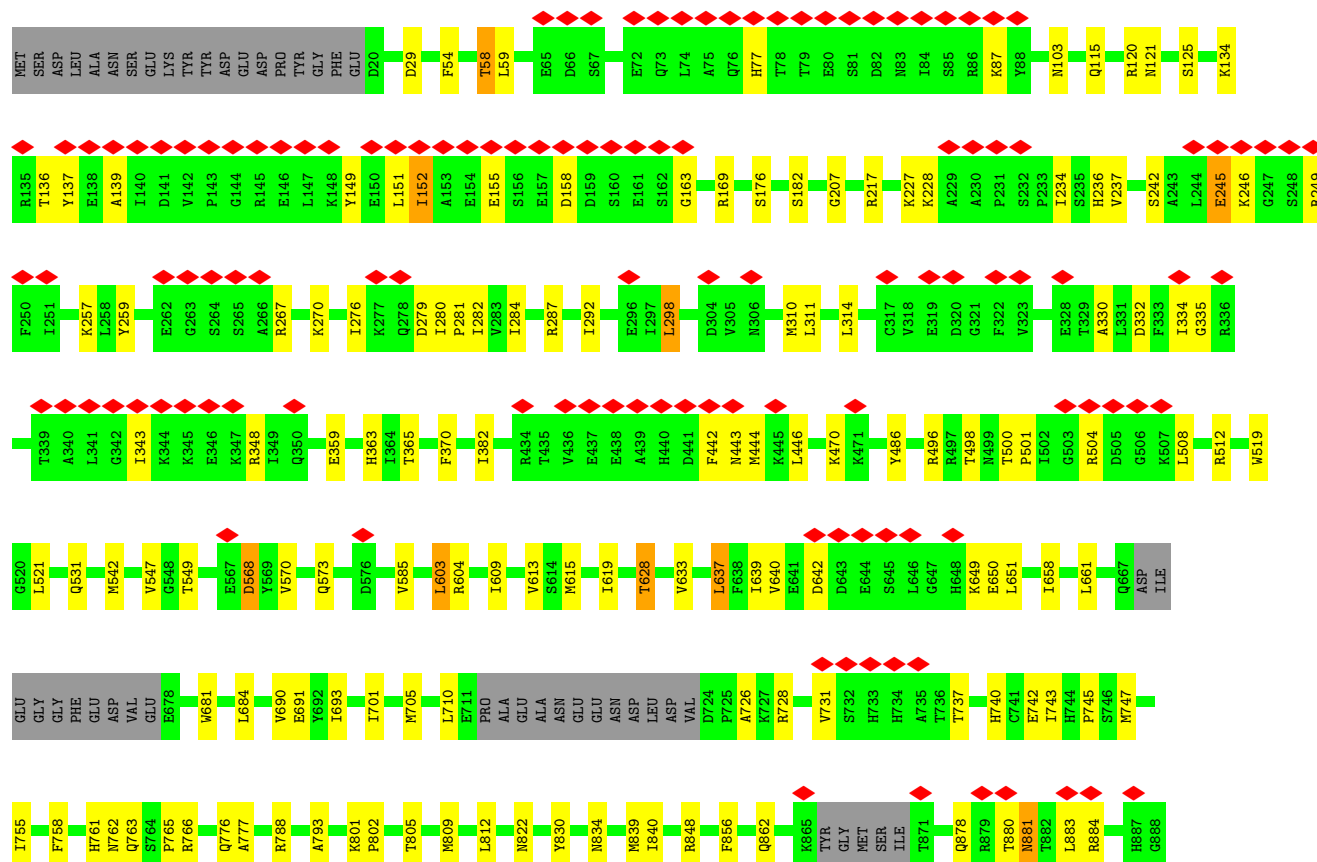
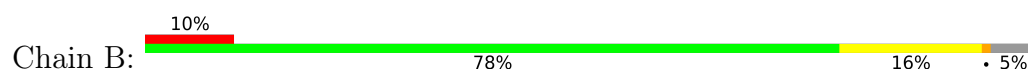
3 Residue-property plots

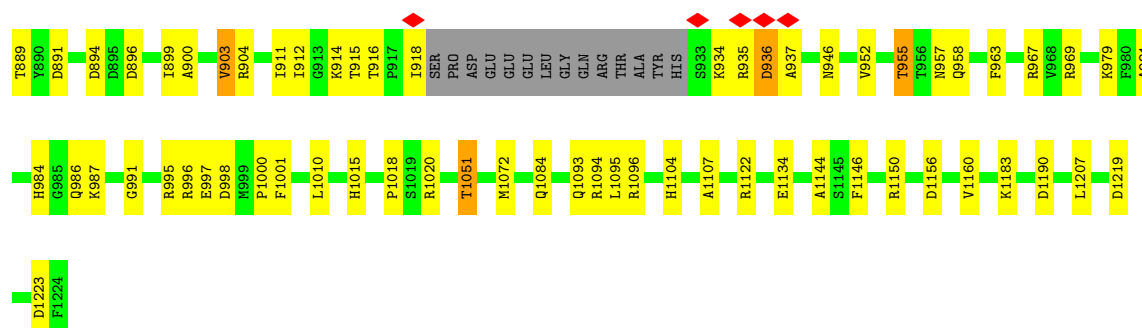
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: DNA-directed RNA polymerase II subunit RPB1

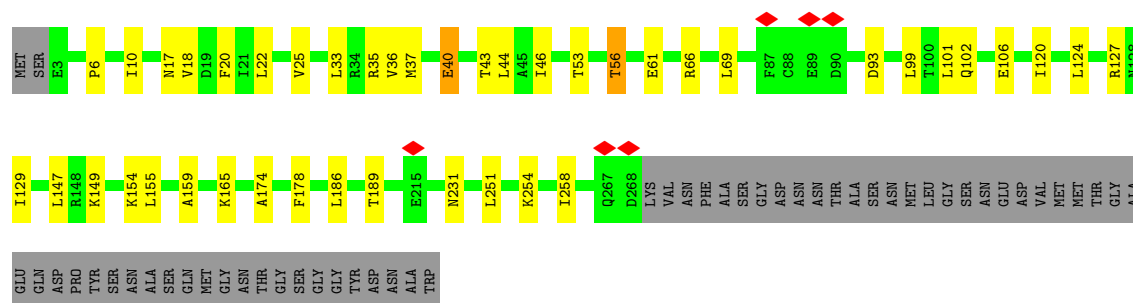


- Molecule 2: DNA-directed RNA polymerase II subunit RPB2

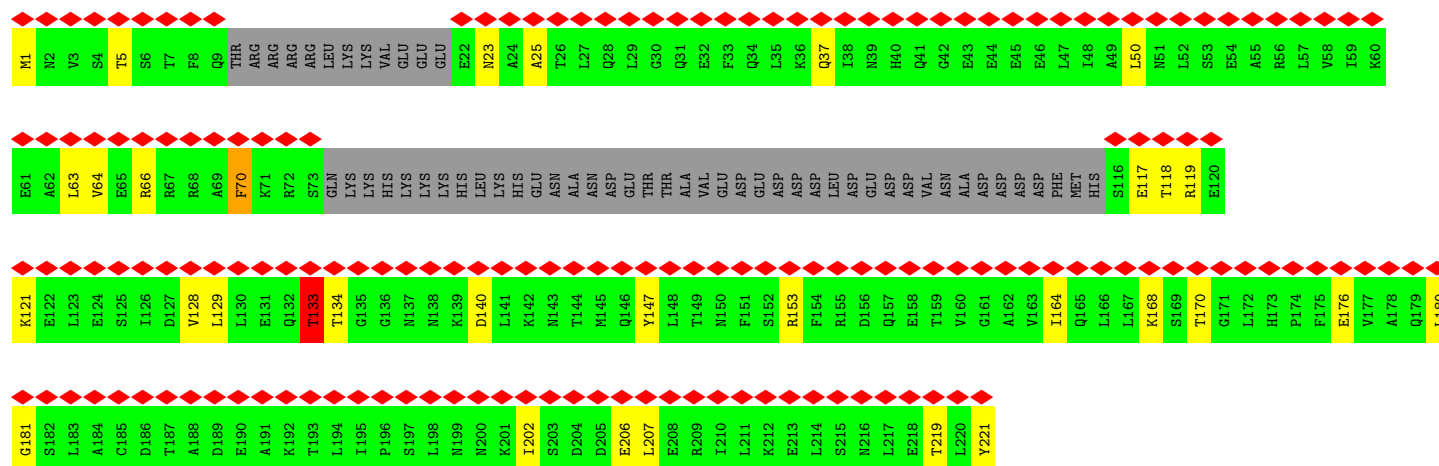
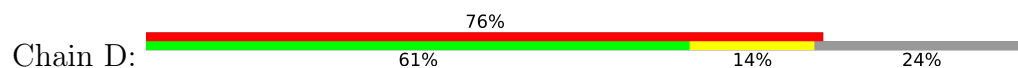




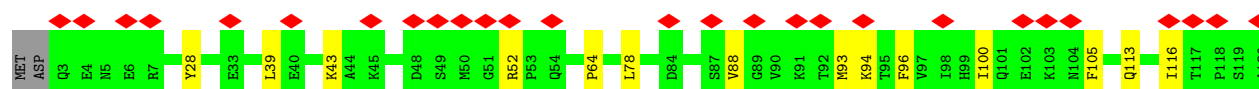
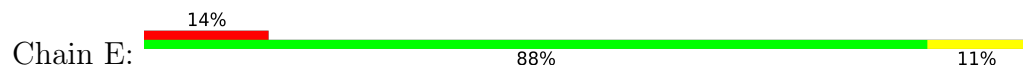
• Molecule 3: DNA-directed RNA polymerase II subunit RPB3



• Molecule 4: DNA-directed RNA polymerase II subunit RPB4



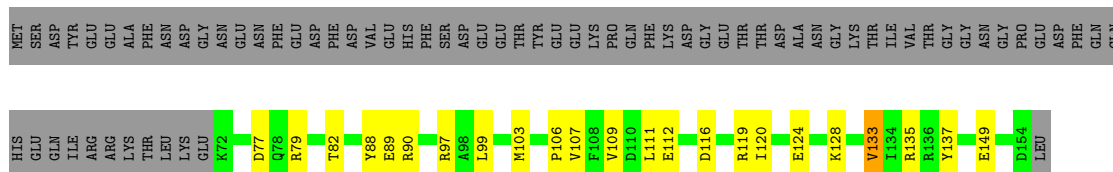
• Molecule 5: DNA-directed RNA polymerases I, II, and III subunit RPABC1





- Molecule 6: DNA-directed RNA polymerases I, II, and III subunit RPABC2

Chain F: 39% 14% 46%



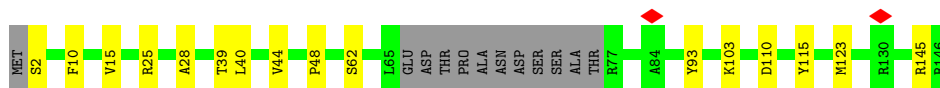
- Molecule 7: DNA-directed RNA polymerase II subunit RPB7

Chain G: 74% 92% 25%



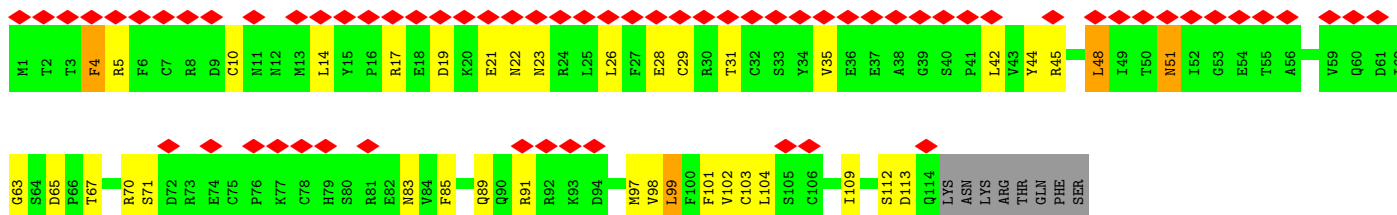
- Molecule 8: DNA-directed RNA polymerases I, II, and III subunit RPABC3

Chain H: 81% 11% 8%



- Molecule 9: DNA-directed RNA polymerase II subunit RPB9

Chain I: 55% 62% 28% 7%



- Molecule 10: DNA-directed RNA polymerases I, II, and III subunit RPABC5

Chain J: 79% 13% 7%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	743682	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	2.489	Depositor
Minimum map value	-1.476	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.052	Depositor
Recommended contour level	0.22	Depositor
Map size ($\text{\AA}$)	346.5, 346.5, 346.5	wwPDB
Map dimensions	420, 420, 420	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.825, 0.825, 0.825	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, MG, DPO, K

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	0.11	0/11193	0.29	0/15132
2	B	0.11	0/9448	0.28	0/12738
3	C	0.12	0/2133	0.28	0/2891
4	D	2.19	8/1336 (0.6%)	0.66	7/1794 (0.4%)
5	E	0.10	0/1780	0.26	0/2395
6	F	0.10	0/682	0.25	0/922
7	G	0.08	0/1368	0.23	0/1844
8	H	0.09	0/1095	0.24	0/1482
9	I	0.10	0/944	0.30	0/1272
10	J	0.11	0/541	0.30	0/727
11	K	0.10	0/922	0.24	0/1244
12	L	0.11	0/361	0.30	0/478
13	N	0.18	0/376	0.39	0/579
14	R	0.14	0/277	0.35	0/432
15	T	0.21	0/596	0.44	0/914
All	All	0.45	8/33052 (0.0%)	0.31	7/44844 (0.0%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	70	PHE	CD2-CE2	35.58	2.45	1.38
4	D	133	THR	CA-CB	34.29	2.07	1.53
4	D	70	PHE	CD1-CE1	32.40	2.35	1.38
4	D	70	PHE	CE1-CZ	31.41	2.32	1.38
4	D	70	PHE	CE2-CZ	30.39	2.29	1.38
4	D	70	PHE	CG-CD1	20.86	1.82	1.38
4	D	70	PHE	CG-CD2	20.81	1.82	1.38
4	D	133	THR	CB-CG2	9.37	1.83	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	133	THR	CA-CB-CG2	14.35	134.89	110.50
4	D	133	THR	CA-CB-OG1	11.61	127.02	109.60
4	D	133	THR	OG1-CB-CG2	-9.31	90.69	109.30
4	D	133	THR	N-CA-C	-8.50	101.96	111.14
4	D	133	THR	CB-CA-C	7.81	123.24	110.90
4	D	133	THR	N-CA-CB	6.11	118.93	110.07
4	D	70	PHE	CD1-CG-CD2	5.74	127.22	118.60

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	10999	0	11076	160	0
2	B	9268	0	9287	131	0
3	C	2095	0	2051	25	0
4	D	1327	0	1331	64	0
5	E	1744	0	1772	11	0
6	F	670	0	690	14	0
7	G	1340	0	1357	30	0
8	H	1077	0	1050	8	0
9	I	926	0	879	27	0
10	J	532	0	542	10	0
11	K	904	0	911	7	0
12	L	359	0	381	12	0
13	N	334	0	178	10	0
14	R	246	0	121	8	0
15	T	537	0	306	10	0
16	A	1	0	0	0	0
17	A	2	0	0	0	0
17	B	1	0	0	0	0
17	C	1	0	0	0	0
17	I	2	0	0	0	0
17	J	1	0	0	0	0
17	L	1	0	0	0	0
18	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	B	9	0	0	0	0
20	A	267	0	0	3	0
20	B	232	0	0	4	0
20	C	68	0	0	0	0
20	E	22	0	0	0	0
20	F	15	0	0	0	0
20	H	10	0	0	0	0
20	J	18	0	0	0	0
20	K	30	0	0	0	0
20	L	4	0	0	0	0
20	R	22	0	0	1	0
20	T	24	0	0	0	0
All	All	33089	0	31932	469	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 (469) 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:D:70:PHE:CD2	4:D:70:PHE:CG	1.82	1.66
4:D:70:PHE:CG	4:D:70:PHE:CD1	1.82	1.66
4:D:70:PHE:CZ	4:D:133:THR:HB	1.35	1.59
4:D:133:THR:CB	4:D:133:THR:CG2	1.83	1.50
4:D:70:PHE:CD1	4:D:133:THR:HA	1.53	1.42
4:D:133:THR:CB	4:D:133:THR:CA	2.07	1.30
4:D:70:PHE:CZ	4:D:133:THR:CB	2.22	1.21
4:D:70:PHE:CZ	4:D:70:PHE:CE2	2.29	1.20
4:D:70:PHE:CE1	4:D:133:THR:HB	1.77	1.18
4:D:70:PHE:CE2	4:D:133:THR:CB	2.26	1.18
4:D:70:PHE:CZ	4:D:70:PHE:CE1	2.32	1.16
4:D:70:PHE:CD1	4:D:70:PHE:CE1	2.35	1.14
4:D:70:PHE:CE1	4:D:133:THR:CB	2.31	1.13
4:D:70:PHE:CD2	4:D:133:THR:N	2.17	1.11
4:D:70:PHE:CD1	4:D:133:THR:CA	2.38	1.06
4:D:70:PHE:CD2	4:D:133:THR:CB	2.39	1.05
4:D:70:PHE:CD2	4:D:70:PHE:CE2	2.45	1.05
4:D:133:THR:HB	4:D:133:THR:CA	1.78	1.04
4:D:70:PHE:CE1	4:D:133:THR:CA	2.43	1.01
4:D:70:PHE:CD1	4:D:133:THR:CB	2.46	0.99
4:D:70:PHE:CE2	4:D:133:THR:OG1	2.18	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:70:PHE:CZ	4:D:133:THR:C	2.45	0.93
4:D:70:PHE:CG	4:D:133:THR:CA	2.51	0.93
4:D:70:PHE:CE2	4:D:133:THR:CA	2.50	0.92
4:D:70:PHE:CD2	4:D:133:THR:OG1	2.21	0.92
4:D:70:PHE:CD2	4:D:133:THR:CA	2.52	0.92
4:D:70:PHE:CG	4:D:133:THR:CB	2.53	0.91
4:D:70:PHE:CZ	4:D:133:THR:CA	2.56	0.89
4:D:70:PHE:CD1	4:D:133:THR:CG2	2.62	0.83
4:D:70:PHE:CG	4:D:133:THR:HA	2.15	0.82
4:D:70:PHE:CE1	4:D:133:THR:HA	2.15	0.80
4:D:70:PHE:CE2	4:D:133:THR:HB	2.17	0.77
2:B:903:VAL:HG12	12:L:61:THR:HG21	1.69	0.75
1:A:1376:THR:O	5:E:212:ARG:NH2	2.20	0.74
1:A:1132:LYS:HG2	1:A:1284:MET:HE1	1.70	0.74
4:D:70:PHE:CE1	4:D:133:THR:C	2.66	0.74
4:D:70:PHE:CE1	4:D:133:THR:CG2	2.72	0.72
3:C:10:ILE:HD13	3:C:20:PHE:HB3	1.71	0.72
1:A:1286:LYS:HD2	1:A:1302:PRO:HB2	1.72	0.72
4:D:133:THR:CG2	4:D:133:THR:OG1	2.36	0.72
2:B:134:LYS:HB3	2:B:155:GLU:HB2	1.70	0.71
2:B:242:SER:HG	2:B:363:HIS:HD1	1.36	0.71
2:B:801:LYS:O	10:J:52:THR:OG1	2.08	0.71
1:A:711:ARG:HE	9:I:97:MET:HG2	1.55	0.70
1:A:541:ILE:HD13	1:A:549:MET:HE1	1.74	0.70
2:B:29:ASP:HB3	2:B:658:ILE:HG12	1.74	0.69
2:B:705:MET:HE2	2:B:745:PRO:HB3	1.73	0.69
5:E:152:LYS:HE3	5:E:154:ILE:HD11	1.74	0.69
1:A:455:MET:HE1	2:B:1134:GLU:HA	1.73	0.69
1:A:968:GLN:HG2	1:A:973:ILE:HD11	1.75	0.69
2:B:642:ASP:HB3	2:B:649:LYS:HG3	1.73	0.69
8:H:44:VAL:HG13	8:H:48:PRO:HA	1.75	0.68
1:A:326:ARG:HG3	1:A:1406:VAL:HG21	1.75	0.68
2:B:762:ASN:HD21	2:B:984:HIS:HD2	1.40	0.68
2:B:987:LYS:NZ	14:R:18:A:OP1	2.23	0.68
2:B:234:ILE:HD13	2:B:257:LYS:HD2	1.76	0.67
2:B:542:MET:HE2	2:B:747:MET:HG3	1.75	0.67
4:D:70:PHE:CE1	4:D:133:THR:HG22	2.30	0.67
7:G:139:ILE:HG22	7:G:140:LYS:HG3	1.77	0.67
4:D:153:ARG:NH1	4:D:180:LEU:O	2.28	0.67
1:A:1267:MET:HA	1:A:1271:ILE:HD13	1.77	0.67
2:B:287:ARG:NH1	2:B:292:ILE:O	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:469:ARG:NH2	2:B:991:GLY:O	2.24	0.66
2:B:889:THR:HG22	2:B:891:ASP:H	1.59	0.66
6:F:111:LEU:HD22	6:F:120:ILE:HD12	1.77	0.66
2:B:998:ASP:OD1	3:C:35:ARG:NH2	2.29	0.66
1:A:1444:MET:HB2	6:F:133:VAL:HG13	1.77	0.65
4:D:70:PHE:CE2	4:D:133:THR:N	2.64	0.65
2:B:862:GLN:O	2:B:914:LYS:NZ	2.28	0.65
2:B:916:THR:HG22	2:B:936:ASP:HB3	1.78	0.65
1:A:147:VAL:HG12	1:A:149:GLU:H	1.62	0.64
2:B:87:LYS:HB3	2:B:137:TYR:HB2	1.78	0.64
2:B:995:ARG:NH1	2:B:997:GLU:OE2	2.30	0.64
7:G:91:VAL:HA	7:G:101:VAL:HA	1.79	0.64
2:B:281:PRO:HD2	2:B:284:ILE:HD12	1.80	0.64
2:B:332:ASP:OD1	2:B:348:ARG:NH1	2.30	0.63
2:B:878:GLN:HB2	2:B:881:ASN:HD21	1.63	0.63
7:G:102:GLN:OE1	7:G:107:LYS:NZ	2.27	0.63
1:A:1308:THR:HG22	1:A:1310:GLY:H	1.63	0.62
7:G:102:GLN:NE2	7:G:104:GLY:O	2.32	0.62
1:A:406:ILE:HB	1:A:431:LYS:HB2	1.81	0.62
4:D:70:PHE:CD1	4:D:133:THR:HG23	2.34	0.62
7:G:89:GLY:HA3	7:G:103:VAL:HG22	1.82	0.62
1:A:120:GLU:OE1	1:A:123:ARG:NH1	2.32	0.62
1:A:789:LYS:HG3	9:I:67:THR:HG22	1.80	0.62
2:B:125:SER:OG	2:B:169:ARG:NH1	2.33	0.62
1:A:875:ALA:HA	1:A:878:ILE:HD12	1.82	0.61
1:A:1009:ASN:OD1	1:A:1012:ARG:NH1	2.33	0.61
2:B:158:ASP:HA	2:B:163:GLY:HA3	1.82	0.61
2:B:443:ASN:HD22	2:B:446:LEU:HG	1.64	0.61
1:A:1283:VAL:HG22	1:A:1307:GLU:HB2	1.82	0.61
1:A:243:PRO:HB2	1:A:245:PRO:HD2	1.83	0.61
1:A:1116:LEU:HD23	1:A:1329:THR:HB	1.83	0.61
1:A:1161:THR:HB	1:A:1170:ILE:HG13	1.81	0.61
7:G:116:PRO:HB2	7:G:119:LEU:HD13	1.83	0.61
2:B:276:ILE:HG21	2:B:280:ILE:HD11	1.83	0.61
2:B:604:ARG:HE	2:B:615:MET:HE2	1.66	0.60
4:D:70:PHE:CG	4:D:133:THR:N	2.69	0.60
1:A:1345:ARG:HG2	1:A:1372:VAL:HG12	1.84	0.60
2:B:137:TYR:HB3	2:B:149:TYR:HB3	1.84	0.59
2:B:848:ARG:NH1	10:J:8:PHE:O	2.34	0.59
1:A:545:GLN:HG2	1:A:549:MET:HE2	1.84	0.59
14:R:9:G:H2'	14:R:10:A:C8	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:T:26:DC:H2'	15:T:27:DT:H71	1.84	0.59
4:D:153:ARG:HH11	4:D:181:GLY:HA2	1.66	0.59
4:D:117:GLU:HB2	4:D:121:LYS:HD2	1.85	0.59
2:B:763:GLN:HG2	2:B:765:PRO:HD2	1.85	0.59
1:A:1111:MET:HE1	1:A:1331:SER:HA	1.85	0.59
2:B:619:ILE:HG13	9:I:65:ASP:HB2	1.84	0.59
2:B:918:ILE:HB	2:B:935:ARG:HB2	1.84	0.59
5:E:88:VAL:HB	5:E:116:ILE:HG12	1.84	0.59
1:A:503:GLN:OE1	6:F:90:ARG:NH2	2.37	0.58
1:A:1100:ARG:HH12	1:A:1111:MET:HE2	1.68	0.58
4:D:66:ARG:HH21	7:G:31:LEU:HD13	1.67	0.58
9:I:21:GLU:OE1	9:I:22:ASN:ND2	2.37	0.58
1:A:1207:LEU:HD21	1:A:1274:ARG:HE	1.68	0.58
1:A:1124:HIS:O	1:A:1130:GLN:NE2	2.36	0.57
3:C:93:ASP:O	3:C:127:ARG:NH2	2.37	0.57
2:B:918:ILE:HD12	2:B:935:ARG:HD2	1.86	0.57
1:A:908:LEU:HD11	1:A:919:ILE:HG21	1.86	0.57
2:B:883:LEU:HB3	2:B:935:ARG:HA	1.87	0.57
9:I:51:ASN:OD1	9:I:51:ASN:N	2.37	0.57
1:A:1129:GLU:OE1	1:A:1132:LYS:NZ	2.35	0.57
1:A:1196:GLU:OE2	9:I:45:ARG:NH2	2.37	0.57
3:C:17:ASN:HD22	3:C:231:ASN:HD21	1.52	0.56
4:D:133:THR:CB	4:D:133:THR:HA	2.25	0.56
3:C:35:ARG:NH1	11:K:41:THR:OG1	2.38	0.56
7:G:147:ILE:HG22	7:G:149:GLY:H	1.69	0.56
4:D:37:GLN:OE1	7:G:5:LYS:NZ	2.38	0.56
2:B:120:ARG:HG2	2:B:955:THR:HG21	1.88	0.56
2:B:249:ARG:NH2	2:B:359:GLU:OE2	2.37	0.56
2:B:896:ASP:OD2	12:L:29:TYR:OH	2.23	0.56
1:A:148:CYS:HB3	1:A:168:GLY:H	1.70	0.56
1:A:18:GLN:HG2	1:A:1418:LEU:HD13	1.88	0.56
2:B:139:ALA:HB2	2:B:149:TYR:HA	1.87	0.56
9:I:83:ASN:HB3	9:I:103:CYS:HA	1.87	0.56
1:A:675:THR:O	1:A:679:ILE:HG12	2.06	0.55
5:E:100:ILE:HG23	5:E:105:PHE:HB2	1.87	0.55
13:N:70:DG:H2''	13:N:71:DA:C8	2.41	0.55
15:T:16:DG:H2''	15:T:17:DG:H5'	1.87	0.55
2:B:996:ARG:NH2	3:C:174:ALA:O	2.39	0.55
4:D:153:ARG:O	4:D:219:THR:OG1	2.21	0.55
3:C:22:LEU:HD11	11:K:101:LEU:HD21	1.89	0.55
7:G:115:MET:HG2	7:G:163:ILE:HD11	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:42:LEU:HD11	9:I:45:ARG:HB2	1.88	0.55
1:A:1197:LEU:HD12	1:A:1209:MET:HE1	1.88	0.55
2:B:822:ASN:O	10:J:48:ARG:NH1	2.34	0.55
3:C:6:PRO:O	11:K:104:ASN:ND2	2.36	0.55
2:B:1094:ARG:NH2	20:B:1424:HOH:O	2.40	0.55
4:D:63:LEU:HB3	4:D:129:LEU:HB2	1.89	0.54
7:G:131:GLN:HG2	7:G:136:VAL:HG13	1.89	0.54
12:L:26:THR:OG1	12:L:40:LEU:O	2.25	0.54
14:R:9:G:O2'	14:R:10:A:OP1	2.24	0.54
14:R:13:G:H2'	14:R:14:G:C8	2.42	0.54
1:A:584:ASN:ND2	20:A:1902:HOH:O	2.31	0.54
1:A:1397:LEU:HB2	1:A:1426:GLU:HG3	1.89	0.54
2:B:762:ASN:HD21	2:B:984:HIS:CD2	2.23	0.54
1:A:1141:THR:HG23	1:A:1273:LEU:HB2	1.90	0.54
7:G:114:LEU:HA	7:G:164:LYS:HE3	1.89	0.54
1:A:500:GLU:OE1	2:B:1146:PHE:N	2.39	0.54
5:E:93:MET:HG3	5:E:123:LEU:HD23	1.88	0.54
10:J:8:PHE:H	10:J:49:MET:HE3	1.73	0.54
1:A:1120:LEU:HD21	1:A:1131:ALA:HB2	1.90	0.53
2:B:365:THR:HG21	2:B:370:PHE:HB2	1.89	0.53
1:A:1147:THR:HB	9:I:48:LEU:HD11	1.90	0.53
14:R:18:A:H2'	14:R:19:G:O4'	2.09	0.53
1:A:407:ARG:NH1	1:A:409:SER:OG	2.41	0.53
2:B:788:ARG:O	2:B:967:ARG:NH1	2.41	0.53
4:D:70:PHE:CZ	4:D:134:THR:N	2.76	0.53
2:B:568:ASP:OD1	2:B:568:ASP:N	2.40	0.53
2:B:915:THR:HB	2:B:934:LYS:HB3	1.91	0.53
3:C:258:ILE:HG13	11:K:42:LEU:HD21	1.91	0.53
6:F:97:ARG:NH2	6:F:106:PRO:O	2.38	0.53
1:A:1409:LEU:HD13	2:B:1207:LEU:HD21	1.91	0.53
2:B:115:GLN:NE2	20:B:1426:HOH:O	2.41	0.53
1:A:587:HIS:CE1	1:A:969:GLN:HG3	2.44	0.53
1:A:663:SER:OG	2:B:1084:GLN:O	2.26	0.53
1:A:1192:LEU:HD11	1:A:1239:ARG:HB3	1.91	0.53
2:B:900:ALA:HB2	12:L:58:LYS:HZ2	1.74	0.52
2:B:969:ARG:NH1	3:C:61:GLU:OE1	2.42	0.52
1:A:899:VAL:HG22	1:A:1029:ARG:HG2	1.91	0.52
2:B:504:ARG:HG2	2:B:531:GLN:HB2	1.91	0.52
1:A:702:LEU:H	1:A:710:LEU:HD11	1.74	0.52
5:E:39:LEU:HG	5:E:43:LYS:HE2	1.91	0.52
7:G:91:VAL:HG12	7:G:101:VAL:HG22	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1113:THR:O	1:A:1113:THR:OG1	2.23	0.52
15:T:20:DC:H5'	15:T:20:DC:C6	2.44	0.52
2:B:705:MET:HE3	2:B:742:GLU:HG3	1.91	0.52
1:A:1231:ASP:OD1	1:A:1231:ASP:N	2.39	0.52
1:A:445:ASN:HB2	1:A:455:MET:HG3	1.91	0.52
4:D:164:ILE:O	4:D:168:LYS:N	2.40	0.52
5:E:127:ILE:HD11	5:E:132:ILE:HD11	1.91	0.52
2:B:217:ARG:NH2	20:B:1421:HOH:O	2.38	0.51
2:B:862:GLN:HG2	2:B:963:PHE:HB2	1.91	0.51
1:A:1161:THR:OG1	1:A:1239:ARG:NH2	2.44	0.51
2:B:613:VAL:HG22	2:B:628:THR:HG23	1.91	0.51
4:D:70:PHE:CE1	4:D:133:THR:O	2.63	0.51
3:C:40:GLU:OE2	3:C:254:LYS:NZ	2.41	0.51
13:N:68:DG:H2''	13:N:69:DG:C8	2.46	0.51
2:B:504:ARG:HG3	2:B:512:ARG:HH22	1.76	0.51
2:B:134:LYS:O	2:B:155:GLU:N	2.41	0.51
1:A:1153:TYR:HE1	9:I:42:LEU:HD13	1.75	0.51
14:R:19:G:P	20:R:101:HOH:O	2.69	0.51
3:C:46:ILE:HD13	3:C:159:ALA:HB2	1.93	0.50
12:L:32:ALA:HB3	12:L:53:HIS:CE1	2.46	0.50
2:B:121:ASN:HA	2:B:207:GLY:HA3	1.92	0.50
3:C:33:LEU:HG	3:C:37:MET:HE2	1.94	0.50
3:C:106:GLU:O	3:C:149:LYS:NZ	2.42	0.50
9:I:10:CYS:SG	9:I:31:THR:OG1	2.59	0.50
2:B:298:LEU:HD12	2:B:314:LEU:HD13	1.94	0.50
1:A:1120:LEU:HD23	1:A:1130:GLN:HE22	1.76	0.50
8:H:110:ASP:OD1	8:H:110:ASP:N	2.45	0.50
15:T:20:DC:H5'	15:T:20:DC:H6	1.76	0.50
1:A:34:LYS:NZ	1:A:85:ASP:OD2	2.44	0.50
1:A:115:LEU:HB2	1:A:122:MET:HE3	1.94	0.50
1:A:698:GLN:HG2	9:I:99:LEU:HD13	1.93	0.50
2:B:900:ALA:HA	12:L:58:LYS:HD3	1.94	0.50
2:B:519:TRP:HZ2	2:B:705:MET:HE1	1.77	0.49
8:H:10:PHE:HB3	8:H:28:ALA:HB1	1.94	0.49
2:B:637:LEU:HD13	2:B:740:HIS:HB3	1.95	0.49
2:B:904:ARG:NH2	12:L:66:GLN:O	2.45	0.49
1:A:1002:GLY:O	1:A:1008:GLN:NE2	2.33	0.49
2:B:486:TYR:CZ	2:B:1096:ARG:HB3	2.47	0.49
7:G:23:LYS:HG3	7:G:56:ILE:HG21	1.95	0.49
9:I:28:GLU:HB3	9:I:35:VAL:HG22	1.95	0.49
1:A:1227:ILE:HB	1:A:1239:ARG:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:996:ASN:HA	1:A:998:LEU:HG	1.95	0.49
2:B:894:ASP:OD2	12:L:58:LYS:NZ	2.41	0.49
1:A:873:MET:HE1	1:A:1057:VAL:HG22	1.95	0.48
1:A:1062:GLU:OE1	6:F:88:TYR:OH	2.28	0.48
4:D:23:ASN:HD21	4:D:25:ALA:HB3	1.78	0.48
8:H:25:ARG:HD2	8:H:39:THR:HG22	1.95	0.48
10:J:14:VAL:HG21	10:J:49:MET:HG2	1.94	0.48
1:A:1317:MET:HA	1:A:1322:ILE:HD11	1.95	0.48
1:A:687:LYS:HD3	1:A:794:PRO:HG2	1.94	0.48
1:A:1442:ASP:HB2	6:F:135:ARG:HB3	1.95	0.48
3:C:6:PRO:HB2	11:K:101:LEU:HG	1.95	0.48
1:A:91:PHE:HZ	1:A:207:ILE:HD12	1.78	0.48
2:B:755:ILE:HG23	2:B:809:MET:HE2	1.95	0.48
1:A:611:GLN:NE2	20:A:1917:HOH:O	2.38	0.48
2:B:444:MET:HE2	2:B:444:MET:HA	1.95	0.48
9:I:101:PHE:HE1	9:I:112:SER:HB3	1.79	0.48
2:B:884:ARG:O	2:B:937:ALA:HB2	2.14	0.48
2:B:878:GLN:O	2:B:881:ASN:ND2	2.46	0.48
9:I:98:VAL:HG11	9:I:113:ASP:HB2	1.95	0.48
1:A:118:HIS:CD2	1:A:152:VAL:HG11	2.49	0.48
1:A:186:LYS:HA	1:A:196:GLU:HB3	1.95	0.48
2:B:134:LYS:HE3	2:B:442:PHE:HZ	1.79	0.48
2:B:136:THR:HG23	2:B:152:ILE:HG23	1.96	0.47
1:A:1114:PRO:HB2	1:A:1311:VAL:HG23	1.95	0.47
1:A:1224:LEU:HD21	1:A:1240:CYS:HB3	1.96	0.47
2:B:640:VAL:HG22	2:B:651:LEU:HD23	1.96	0.47
4:D:64:VAL:HA	4:D:129:LEU:HD22	1.95	0.47
9:I:102:VAL:HG22	9:I:109:ILE:HG12	1.96	0.47
2:B:802:PRO:HG2	2:B:805:THR:HG22	1.97	0.47
1:A:981:LEU:HD23	1:A:1039:LYS:HD2	1.95	0.47
2:B:834:ASN:HB3	2:B:840:ILE:HG13	1.97	0.47
3:C:102:GLN:HG2	3:C:154:LYS:HG3	1.95	0.47
9:I:63:GLY:O	9:I:70:ARG:NH2	2.47	0.47
2:B:661:LEU:HD11	2:B:684:LEU:HD11	1.97	0.47
1:A:439:ASN:OD1	1:A:459:ARG:NH1	2.45	0.47
2:B:176:SER:O	2:B:182:SER:HB3	2.14	0.47
2:B:259:TYR:O	2:B:267:ARG:HA	2.15	0.47
2:B:830:TYR:CZ	2:B:1000:PRO:HD3	2.50	0.47
2:B:900:ALA:HB3	12:L:61:THR:HG22	1.95	0.47
2:B:957:ASN:OD1	2:B:958:GLN:N	2.41	0.47
13:N:71:DA:H2"	13:N:72:DT:C6	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:PRO:HG2	1:A:1067:LEU:HD21	1.96	0.47
7:G:13:LEU:HD23	7:G:13:LEU:HA	1.79	0.47
9:I:83:ASN:HA	9:I:104:LEU:HG	1.97	0.47
1:A:1411:GLU:OE2	1:A:1415:SER:OG	2.32	0.47
2:B:979:LYS:HG2	2:B:1095:LEU:HD12	1.97	0.47
2:B:1104:HIS:HB2	2:B:1122:ARG:HG3	1.97	0.47
12:L:26:THR:OG1	12:L:27:LEU:N	2.48	0.47
1:A:1100:ARG:NH1	1:A:1111:MET:HE2	2.30	0.47
4:D:117:GLU:OE1	4:D:121:LYS:NZ	2.38	0.47
1:A:686:ALA:HB2	1:A:725:ALA:HB2	1.96	0.46
2:B:245:GLU:HB3	2:B:246:LYS:H	1.61	0.46
1:A:1095:THR:OG1	1:A:1096:SER:N	2.48	0.46
1:A:760:GLN:HG2	1:A:765:VAL:HA	1.98	0.46
6:F:107:VAL:HG12	6:F:109:VAL:H	1.80	0.46
8:H:103:LYS:HB3	8:H:115:TYR:HB2	1.96	0.46
1:A:830:LYS:HB2	1:A:1080:THR:HG21	1.97	0.46
11:K:49:GLU:HG3	11:K:94:ILE:HG13	1.97	0.46
2:B:637:LEU:HA	2:B:743:ILE:HG13	1.96	0.46
1:A:195:ASP:OD1	1:A:195:ASP:N	2.48	0.46
1:A:550:LEU:HD21	1:A:561:PRO:HD2	1.98	0.46
1:A:679:ILE:HG23	1:A:729:ALA:HB1	1.97	0.46
1:A:965:GLN:O	1:A:969:GLN:HG2	2.16	0.46
1:A:164:ARG:O	1:A:164:ARG:NE	2.44	0.46
1:A:492:PRO:HB3	1:A:497:THR:HG22	1.98	0.46
1:A:216:VAL:HA	1:A:219:PHE:CE2	2.51	0.46
9:I:5:ARG:HB2	9:I:14:LEU:HB2	1.98	0.46
2:B:311:LEU:HB3	9:I:4:PHE:CE2	2.51	0.45
2:B:603:LEU:HB3	2:B:609:ILE:HG12	1.97	0.45
2:B:726:ALA:HB1	2:B:1051:THR:HG21	1.98	0.45
7:G:143:ILE:HG23	7:G:169:GLY:HA2	1.97	0.45
1:A:1205:LYS:O	1:A:1274:ARG:NH2	2.49	0.45
1:A:1336:MET:HE3	1:A:1380:GLY:HA2	1.98	0.45
2:B:330:ALA:O	2:B:334:ILE:HG12	2.17	0.45
4:D:50:LEU:HD11	7:G:4:ILE:HD11	1.97	0.45
4:D:147:TYR:OH	7:G:88:ASP:O	2.24	0.45
6:F:116:ASP:O	6:F:120:ILE:HG12	2.16	0.45
8:H:93:TYR:HA	8:H:145:ARG:HB3	1.98	0.45
1:A:626:ASN:ND2	20:A:1937:HOH:O	2.49	0.45
1:A:673:GLY:O	1:A:677:ARG:HG3	2.17	0.45
2:B:521:LEU:HD22	2:B:633:VAL:HG12	1.99	0.45
2:B:911:ILE:HG13	2:B:912:ILE:HG13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:R:13:G:H2'	14:R:14:G:H8	1.79	0.45
1:A:32:VAL:HG12	2:B:1183:LYS:NZ	2.32	0.45
1:A:518:LYS:HB3	1:A:626:ASN:HD22	1.81	0.45
1:A:1442:ASP:OD2	6:F:137:TYR:OH	2.31	0.45
2:B:496:ARG:NH1	20:B:1427:HOH:O	2.42	0.45
2:B:570:VAL:HB	2:B:573:GLN:HB3	1.99	0.45
10:J:48:ARG:O	10:J:52:THR:HG22	2.17	0.45
4:D:219:THR:OG1	4:D:219:THR:O	2.35	0.45
10:J:1:MET:HG2	10:J:60:PHE:HE2	1.82	0.45
13:N:68:DG:C2	15:T:23:DG:C2	3.05	0.45
1:A:378:GLU:OE2	1:A:387:ARG:NH2	2.51	0.44
3:C:66:ARG:NH2	10:J:3:VAL:O	2.36	0.44
1:A:260:ASP:OD1	1:A:261:ASP:N	2.50	0.44
1:A:1152:ILE:HG12	9:I:44:TYR:HB3	1.99	0.44
3:C:99:LEU:HG	3:C:120:ILE:HD13	1.98	0.44
12:L:32:ALA:HB3	12:L:53:HIS:HE1	1.82	0.44
2:B:899:ILE:O	2:B:952:VAL:HG21	2.16	0.44
7:G:90:THR:HG23	7:G:140:LYS:HA	2.00	0.44
1:A:455:MET:HE3	1:A:455:MET:HB2	1.69	0.44
1:A:35:ILE:HG22	1:A:270:LEU:HD11	2.00	0.44
4:D:70:PHE:CD1	4:D:70:PHE:CB	2.83	0.44
6:F:99:LEU:HD11	6:F:103:MET:HE3	2.00	0.44
7:G:116:PRO:HD2	7:G:119:LEU:HD22	1.98	0.44
9:I:14:LEU:HD23	9:I:29:CYS:HB3	1.98	0.44
4:D:5:THR:HG22	7:G:7:LEU:HD22	1.99	0.44
13:N:70:DG:H2''	13:N:71:DA:H8	1.79	0.44
1:A:405:VAL:HB	1:A:413:ILE:HB	2.00	0.44
2:B:512:ARG:NH1	2:B:531:GLN:O	2.36	0.44
2:B:650:GLU:HA	2:B:710:LEU:HD13	1.99	0.44
2:B:310:MET:HE2	2:B:310:MET:HB3	1.89	0.43
2:B:282:ILE:HD12	2:B:382:ILE:HD13	1.99	0.43
14:R:9:G:HO2'	14:R:10:A:P	2.38	0.43
1:A:997:LEU:HB3	1:A:1053:PHE:CE1	2.53	0.43
1:A:1328:TYR:CZ	1:A:1350:LYS:HD3	2.53	0.43
2:B:134:LYS:HE3	2:B:442:PHE:CZ	2.53	0.43
8:H:40:LEU:HD13	8:H:123:MET:HB2	2.00	0.43
13:N:73:DC:H1'	13:N:74:DC:O4'	2.17	0.43
1:A:1128:GLN:C	1:A:1130:GLN:H	2.27	0.43
2:B:363:HIS:HD2	2:B:585:VAL:HG22	1.81	0.43
5:E:96:PHE:CZ	5:E:100:ILE:HD11	2.53	0.43
1:A:1436:ILE:HG22	1:A:1437:GLY:H	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:118:THR:OG1	4:D:119:ARG:N	2.51	0.43
7:G:100:GLU:HB3	7:G:107:LYS:HG2	2.01	0.43
1:A:199:LEU:HD23	1:A:199:LEU:HA	1.91	0.43
1:A:752:LYS:HG2	2:B:1015:HIS:O	2.19	0.43
3:C:44:LEU:HD22	3:C:129:ILE:HG12	2.01	0.43
1:A:931:GLU:O	1:A:934:LYS:HG3	2.19	0.43
1:A:1438:THR:HG22	2:B:1144:ALA:HB3	2.01	0.43
7:G:44:TYR:CZ	7:G:105:PRO:HB2	2.54	0.43
1:A:688:LYS:HD3	1:A:688:LYS:HA	1.83	0.43
1:A:1340:GLY:HA2	5:E:183:PRO:HD2	2.00	0.43
4:D:202:ILE:HG23	4:D:207:LEU:HB2	2.01	0.43
11:K:45:LEU:HG	11:K:94:ILE:HD13	2.01	0.43
1:A:1441:PHE:CZ	6:F:89:GLU:HA	2.54	0.43
6:F:109:VAL:HG21	6:F:124:GLU:HA	2.01	0.43
1:A:1400:CYS:HB2	1:A:1405:THR:HG23	2.01	0.42
2:B:793:ALA:HB3	2:B:856:PHE:HB2	2.01	0.42
6:F:116:ASP:HB3	6:F:119:ARG:HB2	2.01	0.42
9:I:19:ASP:O	9:I:23:ASN:N	2.52	0.42
1:A:141:LEU:HD12	1:A:141:LEU:HA	1.83	0.42
1:A:956:LEU:HD21	1:A:1017:LEU:HD23	2.01	0.42
4:D:202:ILE:HD11	4:D:206:GLU:HG3	2.01	0.42
7:G:13:LEU:HG	7:G:26:LEU:HD21	2.01	0.42
15:T:33:DT:H2'	15:T:34:DT:C6	2.55	0.42
1:A:118:HIS:HD2	1:A:152:VAL:HG11	1.83	0.42
1:A:443:LEU:HD12	2:B:1146:PHE:CZ	2.55	0.42
2:B:470:LYS:H	2:B:470:LYS:HD3	1.84	0.42
2:B:500:THR:HA	2:B:501:PRO:HD3	1.84	0.42
3:C:56:THR:HG23	3:C:147:LEU:HD23	2.02	0.42
13:N:61:DG:H2''	13:N:62:DC:C6	2.54	0.42
1:A:1436:ILE:HG22	1:A:1437:GLY:N	2.35	0.42
2:B:639:ILE:HD11	2:B:691:GLU:HG3	2.01	0.42
5:E:94:LYS:HD2	5:E:123:LEU:HD21	2.01	0.42
1:A:553:VAL:HB	1:A:556:TRP:HB2	2.02	0.42
2:B:637:LEU:HD11	2:B:693:ILE:HG13	2.00	0.42
7:G:51:TYR:HA	7:G:54:ILE:HG13	2.02	0.42
15:T:30:DT:H2'	15:T:31:DC:C6	2.55	0.42
1:A:1143:LEU:HD22	1:A:1216:ILE:HD11	2.02	0.42
1:A:1383:SER:OG	1:A:1385:THR:OG1	2.37	0.42
8:H:2:SER:HA	8:H:62:SER:HB2	2.01	0.42
1:A:1420:ASP:OD2	1:A:1422:ARG:NE	2.52	0.42
3:C:66:ARG:NH2	10:J:2:ILE:HG23	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:140:ASP:OD1	4:D:140:ASP:N	2.52	0.42
1:A:16:GLU:HB3	1:A:1418:LEU:HD11	2.01	0.42
1:A:446:ARG:HB2	1:A:487:MET:HE3	2.02	0.42
1:A:922:ASP:OD1	1:A:923:LEU:N	2.53	0.42
2:B:54:PHE:HA	2:B:58:THR:HG23	2.02	0.42
2:B:981:ALA:O	2:B:1093:GLN:HG2	2.19	0.42
6:F:128:LYS:NZ	6:F:149:GLU:O	2.51	0.42
1:A:726:ARG:HD3	1:A:766:GLY:HA3	2.02	0.42
1:A:808:LEU:O	2:B:728:ARG:NH1	2.53	0.42
1:A:886:ILE:HG13	1:A:944:ARG:HG2	2.02	0.42
1:A:1116:LEU:HD11	1:A:1312:ASN:H	1.85	0.42
1:A:1138:ILE:HG22	1:A:1279:ILE:HG21	2.02	0.42
1:A:1158:PRO:O	1:A:1241:ARG:NH1	2.53	0.42
3:C:101:LEU:HD23	3:C:155:LEU:HD12	2.01	0.42
7:G:1:MET:HE2	7:G:1:MET:HB3	1.95	0.42
7:G:1:MET:HG2	7:G:2:PHE:H	1.85	0.42
1:A:399:HIS:CE1	1:A:462:VAL:HG21	2.55	0.42
1:A:549:MET:HG3	1:A:655:PHE:HD2	1.85	0.42
1:A:1174:PHE:HA	1:A:1177:LEU:HD12	2.02	0.42
1:A:1446:ASP:OD1	1:A:1449:SER:N	2.48	0.42
2:B:681:TRP:CH2	2:B:690:VAL:HG11	2.55	0.42
9:I:17:ARG:N	9:I:26:LEU:O	2.43	0.42
1:A:169:ASN:OD1	1:A:169:ASN:N	2.32	0.41
1:A:399:HIS:O	1:A:401:GLY:N	2.51	0.41
1:A:919:ILE:HD13	1:A:919:ILE:HA	1.88	0.41
1:A:1194:ARG:HH21	1:A:1237:ILE:HG21	1.85	0.41
2:B:227:LYS:HG2	2:B:236:HIS:CE1	2.55	0.41
1:A:993:LEU:HD22	1:A:1046:LEU:HD22	2.03	0.41
1:A:528:LEU:HD23	1:A:751:SER:HA	2.02	0.41
1:A:839:ARG:NH2	1:A:1402:PHE:O	2.53	0.41
2:B:758:PHE:HB3	2:B:761:HIS:CD2	2.54	0.41
3:C:36:VAL:HG21	3:C:251:LEU:HD13	2.02	0.41
15:T:17:DG:H2"	15:T:18:DA:C8	2.56	0.41
1:A:106:VAL:HG11	1:A:214:ILE:HD12	2.03	0.41
1:A:172:PRO:HB3	1:A:185:TRP:CE2	2.56	0.41
4:D:70:PHE:CD2	4:D:70:PHE:CB	2.83	0.41
7:G:129:SER:HB2	7:G:137:ILE:O	2.20	0.41
1:A:202:LEU:HB3	1:A:207:ILE:HD11	2.02	0.41
1:A:1209:MET:HG3	1:A:1229:SER:O	2.21	0.41
13:N:73:DC:H5'	13:N:73:DC:C6	2.56	0.41
2:B:701:ILE:HB	2:B:740:HIS:HD2	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1001:PHE:HE2	3:C:178:PHE:HB3	1.86	0.41
2:B:1156:ASP:N	2:B:1156:ASP:OD1	2.54	0.41
9:I:85:PHE:HB3	9:I:101:PHE:CD2	2.55	0.41
13:N:71:DA:H2''	13:N:72:DT:C5	2.54	0.41
2:B:335:GLY:HA3	2:B:348:ARG:HB3	2.03	0.41
2:B:737:THR:HG21	9:I:70:ARG:NH1	2.35	0.41
9:I:71:SER:OG	9:I:83:ASN:OD1	2.34	0.41
1:A:3:GLY:O	1:A:75:ASN:ND2	2.50	0.41
1:A:23:SER:OG	1:A:25:GLU:OE1	2.39	0.41
1:A:1279:ILE:HG23	1:A:1308:THR:HG23	2.02	0.41
2:B:777:ALA:HB1	2:B:1093:GLN:HB2	2.03	0.41
5:E:28:TYR:HA	5:E:64:PRO:HA	2.01	0.41
7:G:137:ILE:HG23	7:G:143:ILE:HG13	2.02	0.41
15:T:36:DT:H2'	15:T:37:DC:C6	2.56	0.41
1:A:767:GLN:HA	1:A:799:PHE:HA	2.03	0.41
1:A:1189:SER:HB2	1:A:1191:TRP:HD1	1.86	0.41
1:A:1198:ASP:HB3	1:A:1201:ALA:HB3	2.03	0.40
2:B:137:TYR:HE1	2:B:151:LEU:HD13	1.85	0.40
10:J:1:MET:HG2	10:J:60:PHE:CE2	2.56	0.40
1:A:346:ASP:HB3	2:B:1107:ALA:O	2.21	0.40
1:A:1095:THR:HG23	1:A:1100:ARG:HB2	2.03	0.40
2:B:270:LYS:HB3	2:B:279:ASP:HB3	2.01	0.40
2:B:839:MET:HB2	2:B:1010:LEU:HD11	2.03	0.40
4:D:1:MET:SD	4:D:1:MET:N	2.74	0.40
15:T:28:DG:H2''	15:T:29:DC:C5	2.56	0.40
1:A:915:SER:O	1:A:919:ILE:HG12	2.21	0.40
1:A:1143:LEU:O	1:A:1147:THR:OG1	2.26	0.40
2:B:228:LYS:HB2	2:B:234:ILE:O	2.21	0.40
2:B:776:GLN:HB3	2:B:1096:ARG:HG2	2.02	0.40
7:G:65:ASP:OD1	7:G:65:ASP:N	2.54	0.40
12:L:53:HIS:CE1	12:L:55:ILE:HB	2.55	0.40
1:A:402:ALA:HA	1:A:434:ARG:HA	2.02	0.40
1:A:752:LYS:HG3	2:B:1018:PRO:HG2	2.03	0.40
1:A:1170:ILE:HG23	1:A:1174:PHE:CD2	2.56	0.40
3:C:124:LEU:O	3:C:127:ARG:HG2	2.21	0.40
13:N:63:DA:H1'	13:N:64:DG:H5'	2.03	0.40
1:A:842:VAL:O	1:A:846:GLU:HB3	2.21	0.40
2:B:766:ARG:CZ	2:B:1020:ARG:HG3	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1381/1733 (80%)	1328 (96%)	52 (4%)	1 (0%)	48	57
2	B	1154/1224 (94%)	1125 (98%)	29 (2%)	0	100	100
3	C	264/318 (83%)	262 (99%)	2 (1%)	0	100	100
4	D	161/221 (73%)	157 (98%)	4 (2%)	0	100	100
5	E	211/215 (98%)	201 (95%)	9 (4%)	1 (0%)	24	26
6	F	81/155 (52%)	80 (99%)	1 (1%)	0	100	100
7	G	169/171 (99%)	164 (97%)	5 (3%)	0	100	100
8	H	130/146 (89%)	130 (100%)	0	0	100	100
9	I	112/122 (92%)	106 (95%)	6 (5%)	0	100	100
10	J	63/70 (90%)	63 (100%)	0	0	100	100
11	K	110/120 (92%)	109 (99%)	1 (1%)	0	100	100
12	L	43/70 (61%)	40 (93%)	2 (5%)	1 (2%)	5	3
All	All	3879/4565 (85%)	3765 (97%)	111 (3%)	3 (0%)	49	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	958	VAL
12	L	45	ALA
5	E	52	ARG

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	1223/1520 (80%)	1170 (96%)	53 (4%)	26	33
2	B	1010/1061 (95%)	977 (97%)	33 (3%)	33	43
3	C	234/274 (85%)	224 (96%)	10 (4%)	26	33
4	D	149/200 (74%)	144 (97%)	5 (3%)	32	42
5	E	195/197 (99%)	190 (97%)	5 (3%)	40	53
6	F	73/137 (53%)	68 (93%)	5 (7%)	14	16
7	G	152/152 (100%)	147 (97%)	5 (3%)	33	43
8	H	118/128 (92%)	117 (99%)	1 (1%)	73	82
9	I	108/116 (93%)	102 (94%)	6 (6%)	19	23
10	J	60/65 (92%)	58 (97%)	2 (3%)	33	43
11	K	97/102 (95%)	95 (98%)	2 (2%)	47	60
12	L	40/57 (70%)	40 (100%)	0	100	100
All	All	3459/4009 (86%)	3332 (96%)	127 (4%)	31	39

All (127) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	THR
1	A	22	PHE
1	A	50	ILE
1	A	93	VAL
1	A	141	LEU
1	A	167	CYS
1	A	169	ASN
1	A	196	GLU
1	A	232	GLU
1	A	241	VAL
1	A	266	LEU
1	A	287	HIS
1	A	335	ARG
1	A	351	THR
1	A	379	VAL
1	A	380	VAL
1	A	391	LEU
1	A	443	LEU
1	A	444	PHE
1	A	447	GLN
1	A	451	HIS
1	A	472	LEU

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Mol	Chain	Res	Type
1	A	474	VAL
1	A	501	LEU
1	A	538	ASP
1	A	550	LEU
1	A	618	GLU
1	A	683	ILE
1	A	698	GLN
1	A	703	THR
1	A	822	GLU
1	A	833	GLU
1	A	838	GLN
1	A	890	ASP
1	A	895	LYS
1	A	904	THR
1	A	907	THR
1	A	912	LEU
1	A	924	LYS
1	A	932	GLU
1	A	934	LYS
1	A	977	LYS
1	A	999	VAL
1	A	1067	LEU
1	A	1095	THR
1	A	1113	THR
1	A	1128	GLN
1	A	1169	ILE
1	A	1231	ASP
1	A	1232	ASN
1	A	1366	ARG
1	A	1377	THR
1	A	1438	THR
2	B	58	THR
2	B	59	LEU
2	B	77	HIS
2	B	103	ASN
2	B	152	ILE
2	B	237	VAL
2	B	245	GLU
2	B	298	LEU
2	B	343	ILE
2	B	498	THR
2	B	508	LEU

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Mol	Chain	Res	Type
2	B	547	VAL
2	B	549	THR
2	B	568	ASP
2	B	603	LEU
2	B	628	THR
2	B	637	LEU
2	B	731	VAL
2	B	812	LEU
2	B	880	THR
2	B	881	ASN
2	B	903	VAL
2	B	936	ASP
2	B	946	ASN
2	B	955	THR
2	B	986	GLN
2	B	1051	THR
2	B	1072	MET
2	B	1150	ARG
2	B	1160	VAL
2	B	1190	ASP
2	B	1219	ASP
2	B	1223	ASP
3	C	18	VAL
3	C	25	VAL
3	C	40	GLU
3	C	43	THR
3	C	53	THR
3	C	56	THR
3	C	69	LEU
3	C	165	LYS
3	C	186	LEU
3	C	189	THR
4	D	128	VAL
4	D	133	THR
4	D	170	THR
4	D	176	GLU
4	D	221	TYR
5	E	78	LEU
5	E	113	GLN
5	E	175	LEU
5	E	196	VAL
5	E	200	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
6	F	77	ASP
6	F	79	ARG
6	F	82	THR
6	F	112	GLU
6	F	133	VAL
7	G	50	ASP
7	G	56	ILE
7	G	77	VAL
7	G	91	VAL
7	G	92	VAL
8	H	15	VAL
9	I	4	PHE
9	I	48	LEU
9	I	51	ASN
9	I	89	GLN
9	I	91	ARG
9	I	99	LEU
10	J	1	MET
10	J	36	LEU
11	K	31	VAL
11	K	101	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (56) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	68	GLN
1	A	71	GLN
1	A	118	HIS
1	A	311	GLN
1	A	316	GLN
1	A	390	GLN
1	A	394	ASN
1	A	397	ASN
1	A	425	GLN
1	A	427	GLN
1	A	451	HIS
1	A	479	ASN
1	A	626	ASN
1	A	631	HIS
1	A	877	HIS
1	A	1033	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	1040	GLN
1	A	1052	GLN
1	A	1130	GLN
1	A	1171	GLN
1	A	1188	GLN
1	A	1203	ASN
2	B	76	GLN
2	B	115	GLN
2	B	121	ASN
2	B	215	GLN
2	B	350	GLN
2	B	383	ASN
2	B	395	GLN
2	B	443	ASN
2	B	449	ASN
2	B	499	ASN
2	B	572	HIS
2	B	740	HIS
2	B	843	GLN
2	B	878	GLN
2	B	881	ASN
2	B	984	HIS
2	B	986	GLN
3	C	231	ASN
3	C	264	GLN
4	D	157	GLN
4	D	179	GLN
5	E	8	ASN
5	E	54	GLN
5	E	101	GLN
7	G	57	GLN
7	G	71	ASN
8	H	11	GLN
8	H	33	GLN
8	H	83	GLN
9	I	12	ASN
9	I	89	GLN
9	I	114	GLN
12	L	53	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
14	R	11/19 (57%)	2 (18%)	1 (9%)

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
14	R	10	A
14	R	19	G

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
14	R	9	G

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 11 ligands modelled in this entry, 10 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
19	DPO	B	1301	-	6,8,8	0.73	0	13,13,13	1.12	1 (7%)

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
19	DPO	B	1301	-	-	0/6/6/6	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
19	B	1301	DPO	P2-O4-P1	-2.65	123.75	132.83

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.

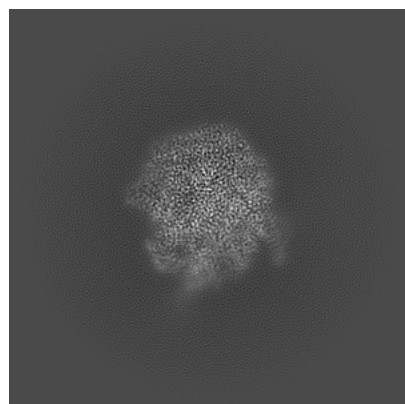
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-54374. These allow visual inspection of the internal detail of the map and identification of artifacts.

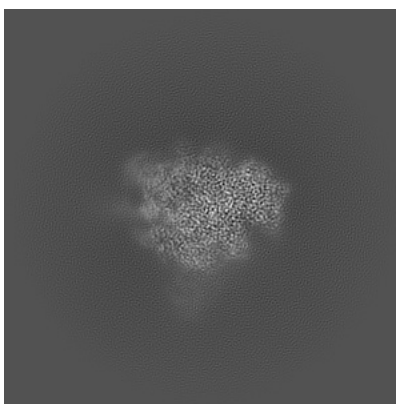
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

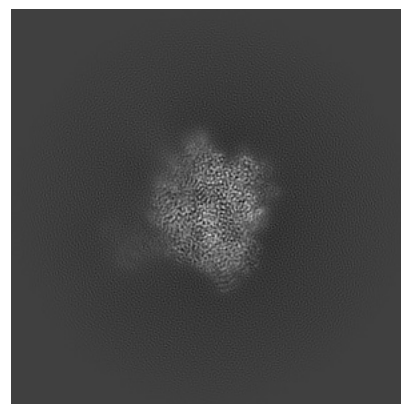
6.1.1 Primary map



X

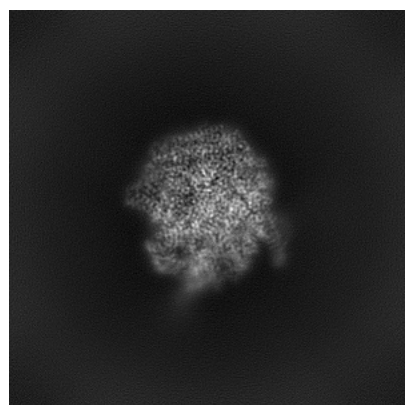


Y

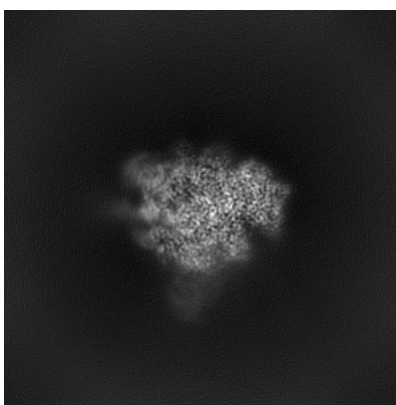


Z

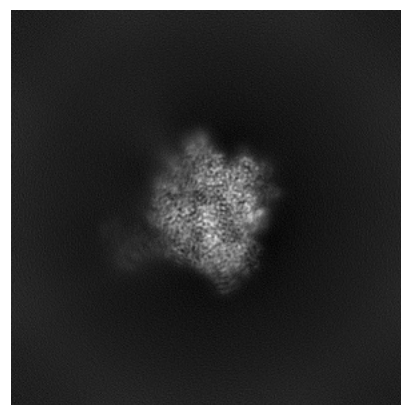
6.1.2 Raw map



X



Y

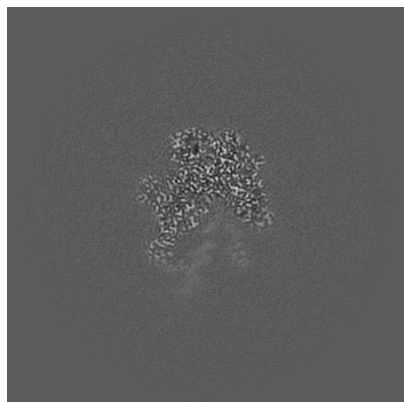


Z

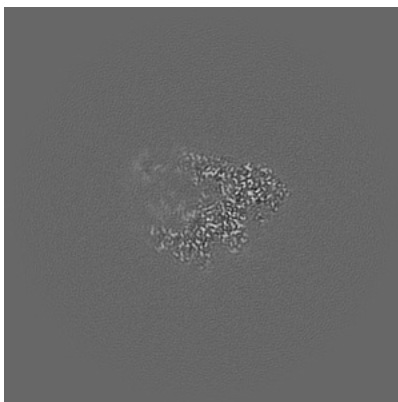
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

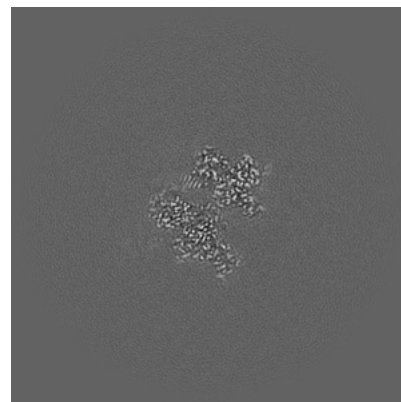
6.2.1 Primary map



X Index: 210

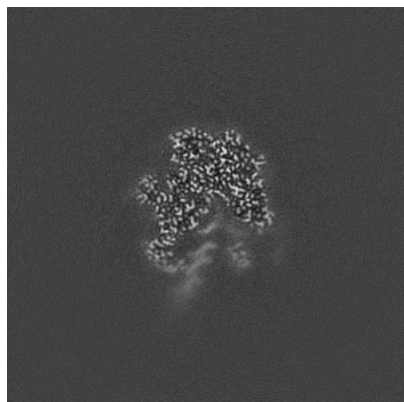


Y Index: 210

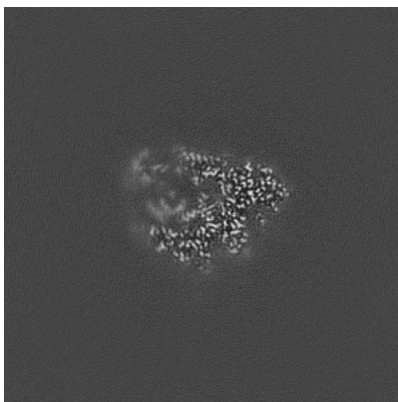


Z Index: 210

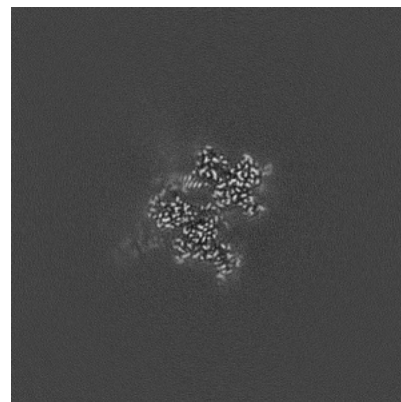
6.2.2 Raw map



X Index: 210



Y Index: 210

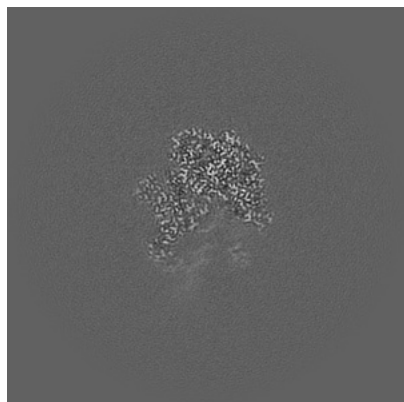


Z Index: 210

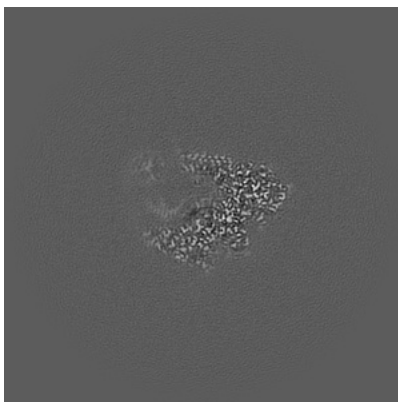
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

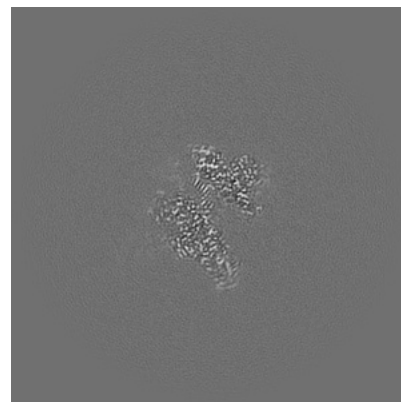
6.3.1 Primary map



X Index: 211

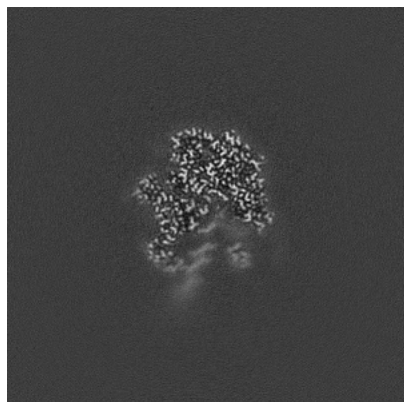


Y Index: 207

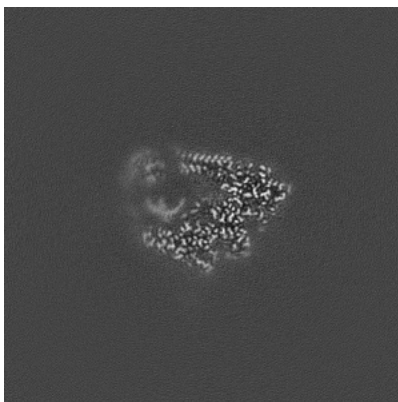


Z Index: 217

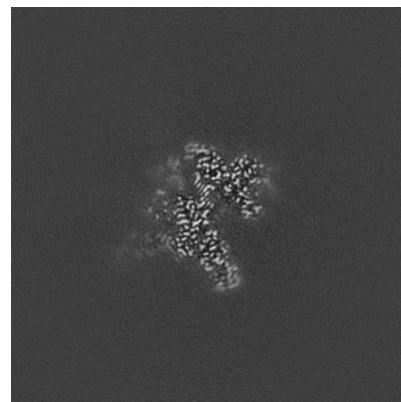
6.3.2 Raw map



X Index: 211



Y Index: 206

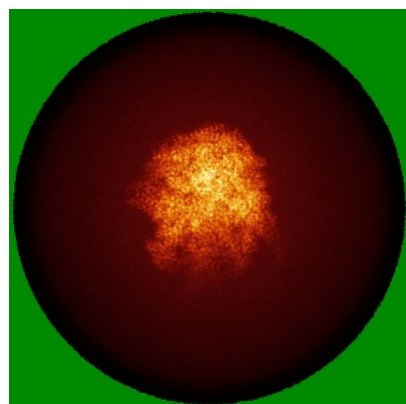


Z Index: 218

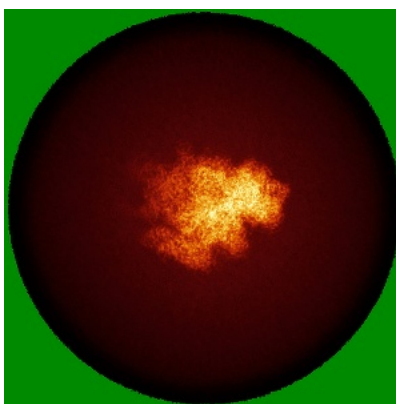
The images above show the largest variance slices of the map in three orthogonal directions.

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

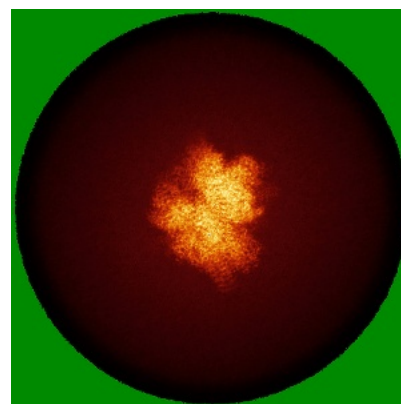
6.4.1 Primary map



X

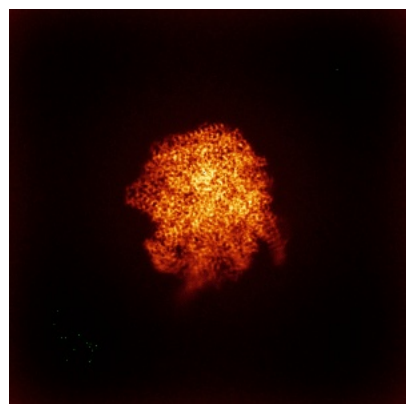


Y

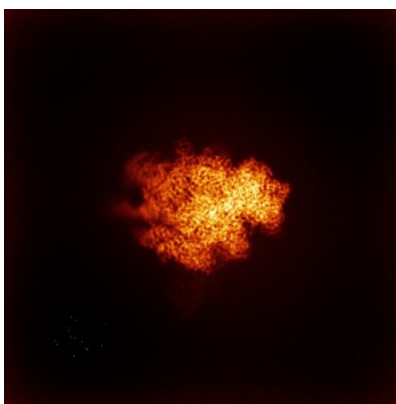


Z

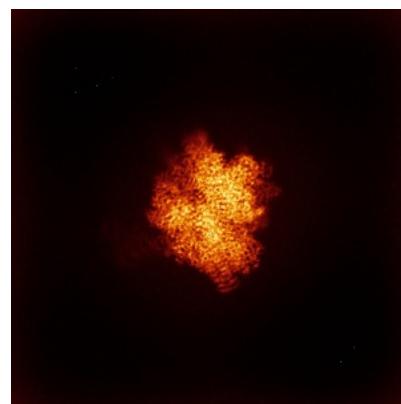
6.4.2 Raw map



X



Y

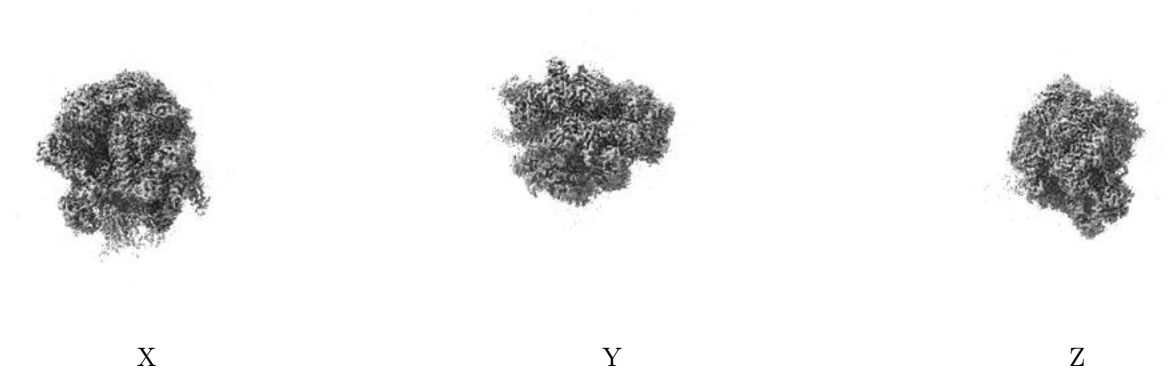


Z

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.

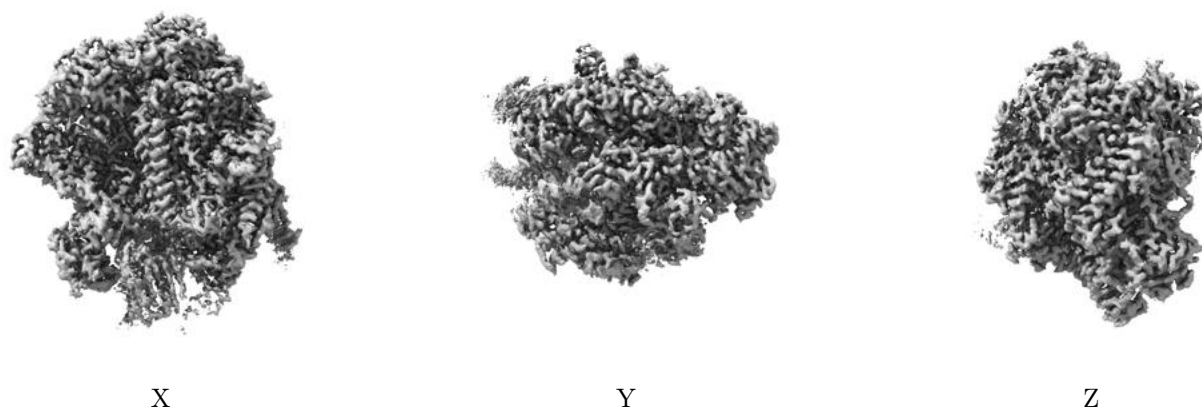
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.22. 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.

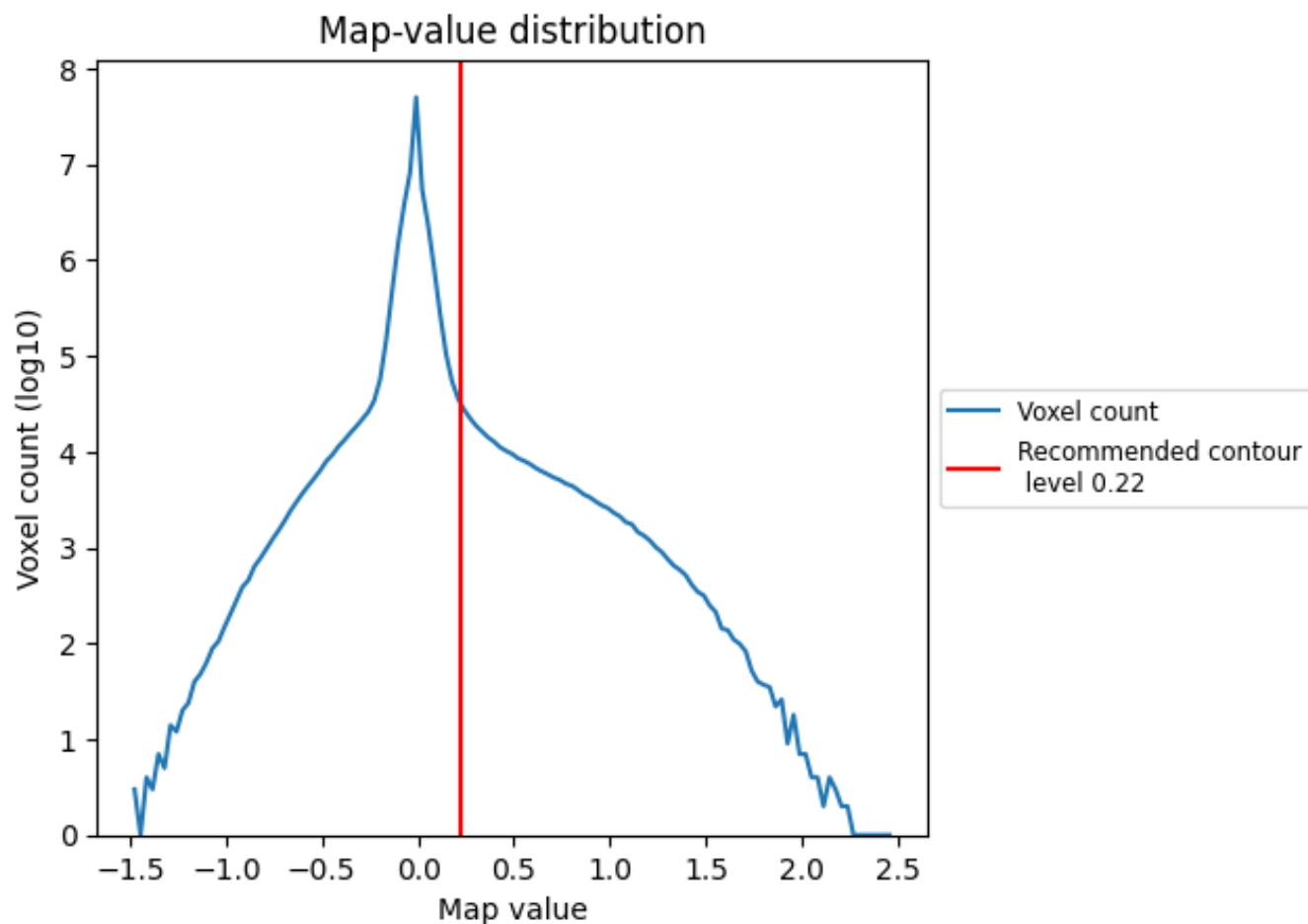
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

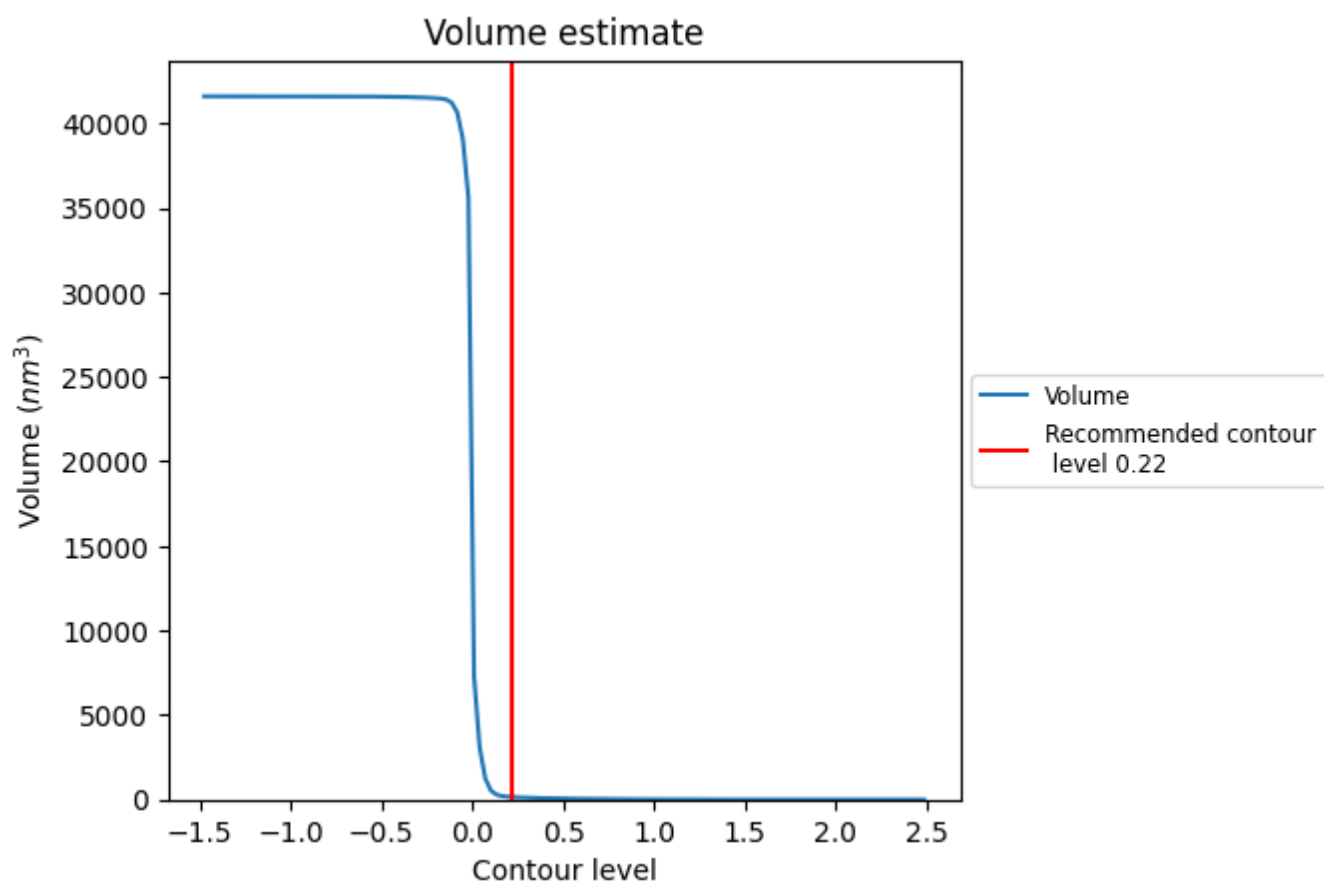
This section contains the results of statistical analysis of the map.

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.

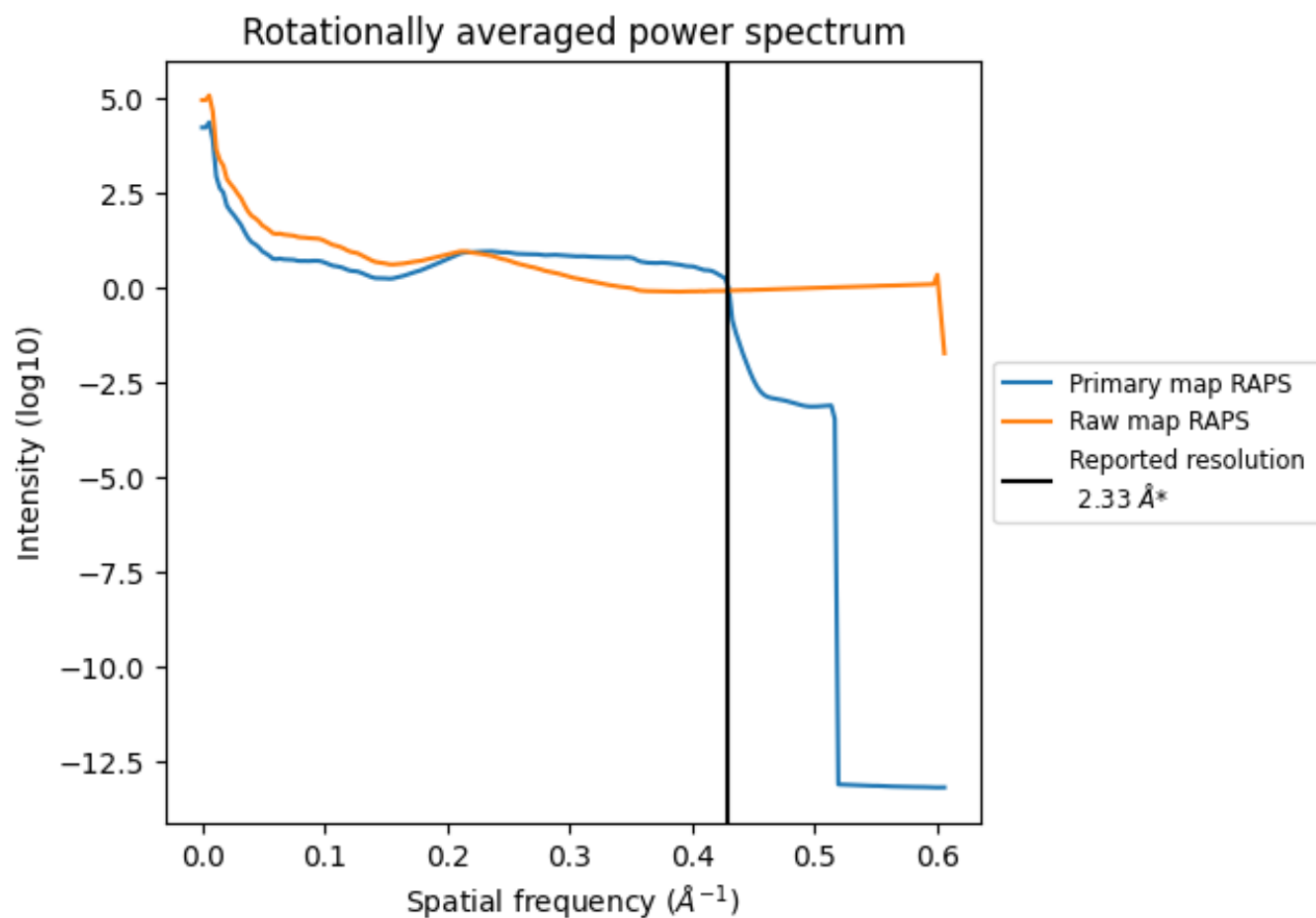
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 150 nm^3 ; this corresponds to an approximate mass of 136 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 ⓘ

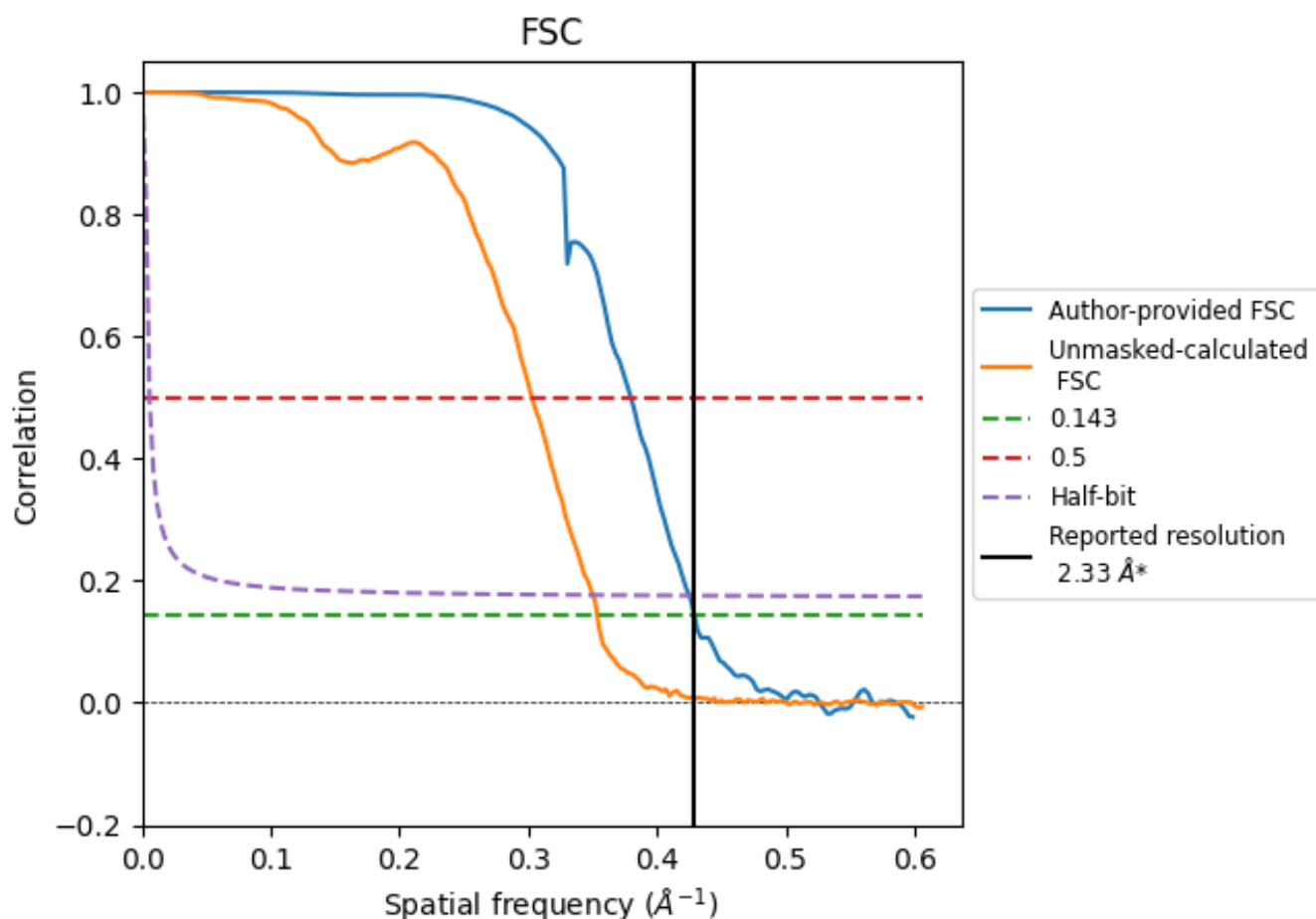


*Reported resolution corresponds to spatial frequency of 0.429 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.429 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

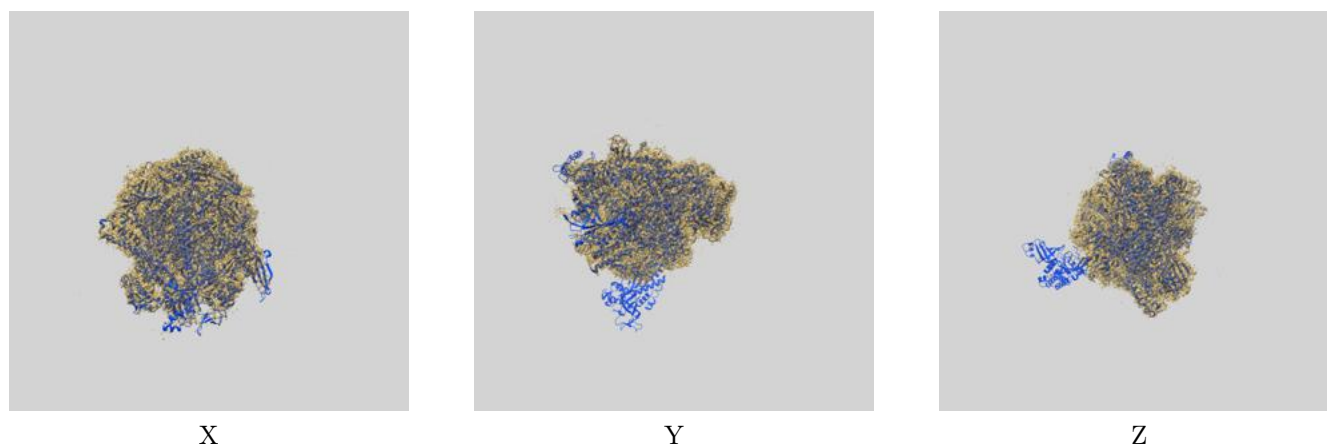
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.33	-	-
Author-provided FSC curve	2.33	2.63	2.35
Unmasked-calculated*	2.83	3.30	2.86

*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 2.83 differs from the reported value 2.33 by more than 10 %

9 Map-model fit [i](#)

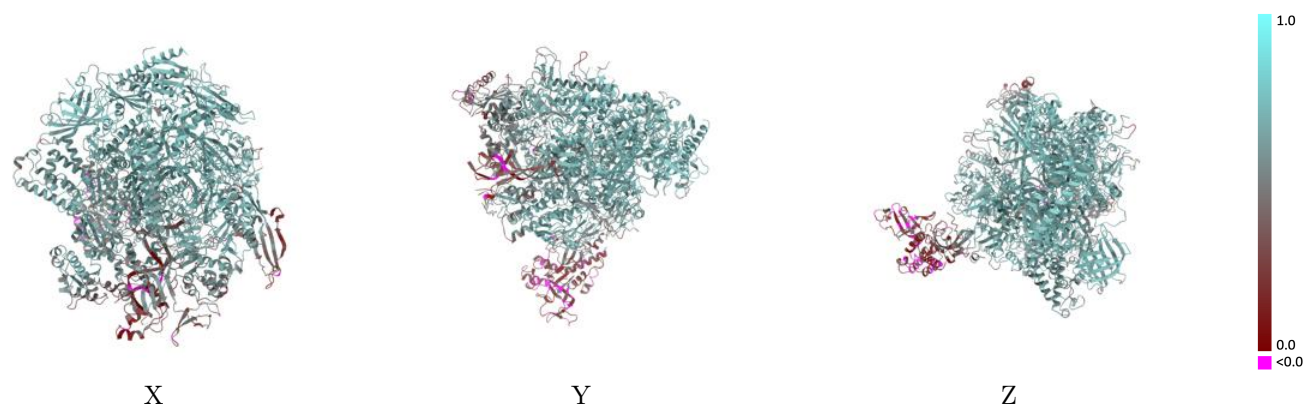
This section contains information regarding the fit between EMDB map EMD-54374 and PDB model 9RYB. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



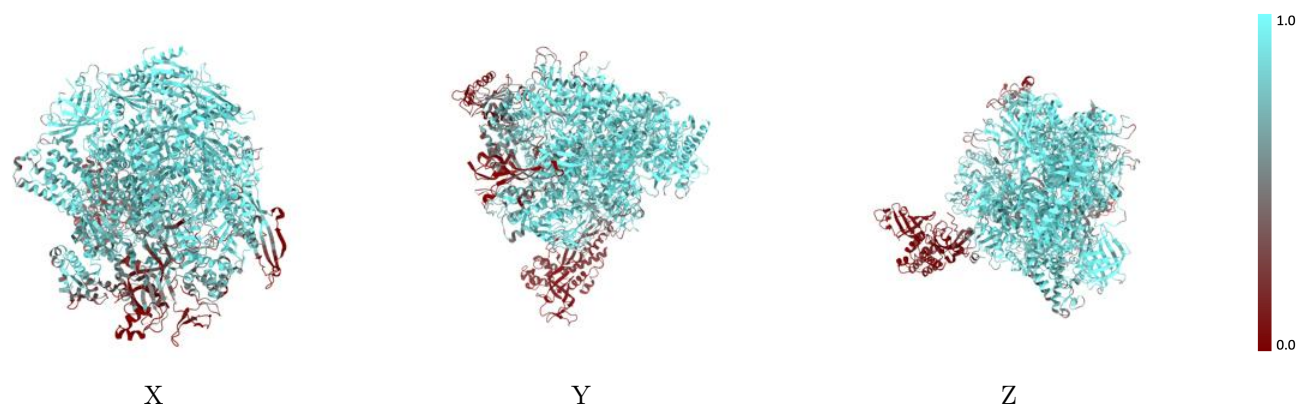
The images above show the 3D surface view of the map at the recommended contour level 0.22 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



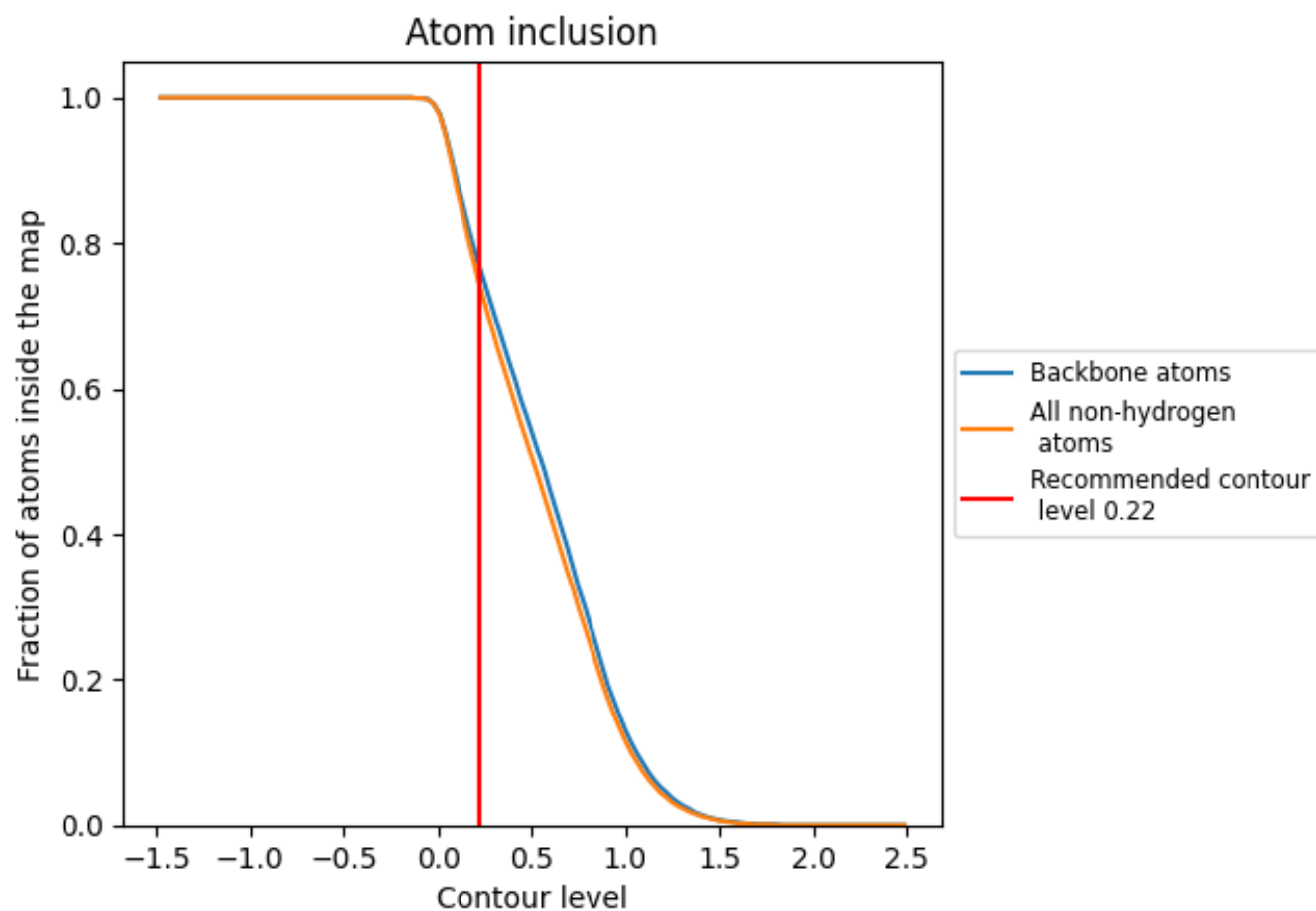
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.22).

9.4 Atom inclusion [i](#)



At the recommended contour level, 77% of all backbone atoms, 75% 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 (0.22) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7460	0.5860
A	0.8210	0.6240
B	0.8250	0.6260
C	0.9150	0.6710
D	0.0020	0.1590
E	0.7680	0.5980
F	0.9190	0.6620
G	0.0830	0.2980
H	0.8850	0.6520
I	0.3990	0.4870
J	0.9500	0.6850
K	0.9350	0.6870
L	0.6510	0.5630
N	0.2390	0.2060
R	0.9020	0.6250
T	0.5810	0.4530

1.0

0.0

<0.0