



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 05:40 PM UTC

PDB ID : 5IP9 / pdb\_00005ip9  
Title : Structure of RNA Polymerase II-TFIIF complex  
Authors : Plaschka, C.; Hantsche, M.; Dienemann, C.; Burzinski, C.; Plitzko, J.; Cramer, P.  
Deposited on : 2016-03-09  
Resolution : 3.90 Å (reported)

This is a wwPDB X-ray Structure 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/XrayValidationReportHelp>

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:

MolProbity : 4-5-2 with Phenix2.0  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

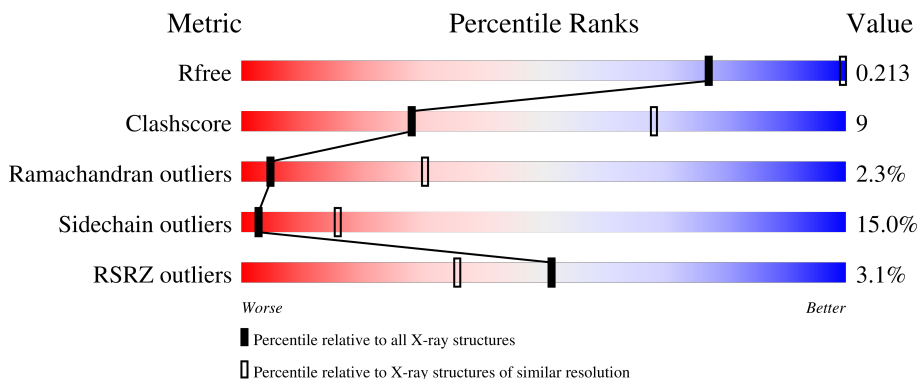
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

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)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1270 (4.10-3.70)
Clashscore	190562	1034 (4.08-3.72)
Ramachandran outliers	187476	1251 (4.10-3.70)
Sidechain outliers	187428	1243 (4.10-3.70)
RSRZ outliers	180081	1269 (4.10-3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1732	 2% 52% 24% 5% 18%
2	B	1223	 4% 60% 26% 5% 9%
3	C	266	 0% 66% 28% 6%
4	D	221	 4% 56% 18% 8% 19%
5	E	214	 0% 76% 22% 0%

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Mol	Chain	Length	Quality of chain
6	F	87	<p>2% 69% 26% 5%</p>
7	G	171	<p>1% 67% 27% 6%</p>
8	H	145	<p>6% 56% 30% 5% 8%</p>
9	I	119	<p>8% 68% 26% 5%</p>
10	J	65	<p>3% 54% 37% 9%</p>
11	K	115	<p>1% 72% 25% 2%</p>
12	L	46	<p>15% 50% 39% 11%</p>
13	Q	15	<p>47% 80% 13% 7%</p>

## 2 Entry composition [i](#)

There are 15 unique types of molecules in this entry. The entry contains 31339 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1421	11190	7052	1957	2119	62	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	1118	8876	5620	1557	1644	55	0	0	0

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	180	1440	890	256	291	3	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	E	214	1752	1111	309	321	11	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
6	F	87	705	451	119	132	3	0	0	0

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
8	H	133	1068	673	180	211	4	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
9	I	119	971	596	179	186	10	0	0	0

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
11	K	115	920	590	157	171	2	0	0	1

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
12	L	46	363	224	72	63	4	0	0	0

- Molecule 13 is a protein called PHE-ILE-LYS-ARG-ASP-ARG-MET-ARG-ARG-ASN-PHE-LEU-ARG-MET-ARG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
13	Q	15	78	47	15	15	1	0	0	0

- Molecule 14 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

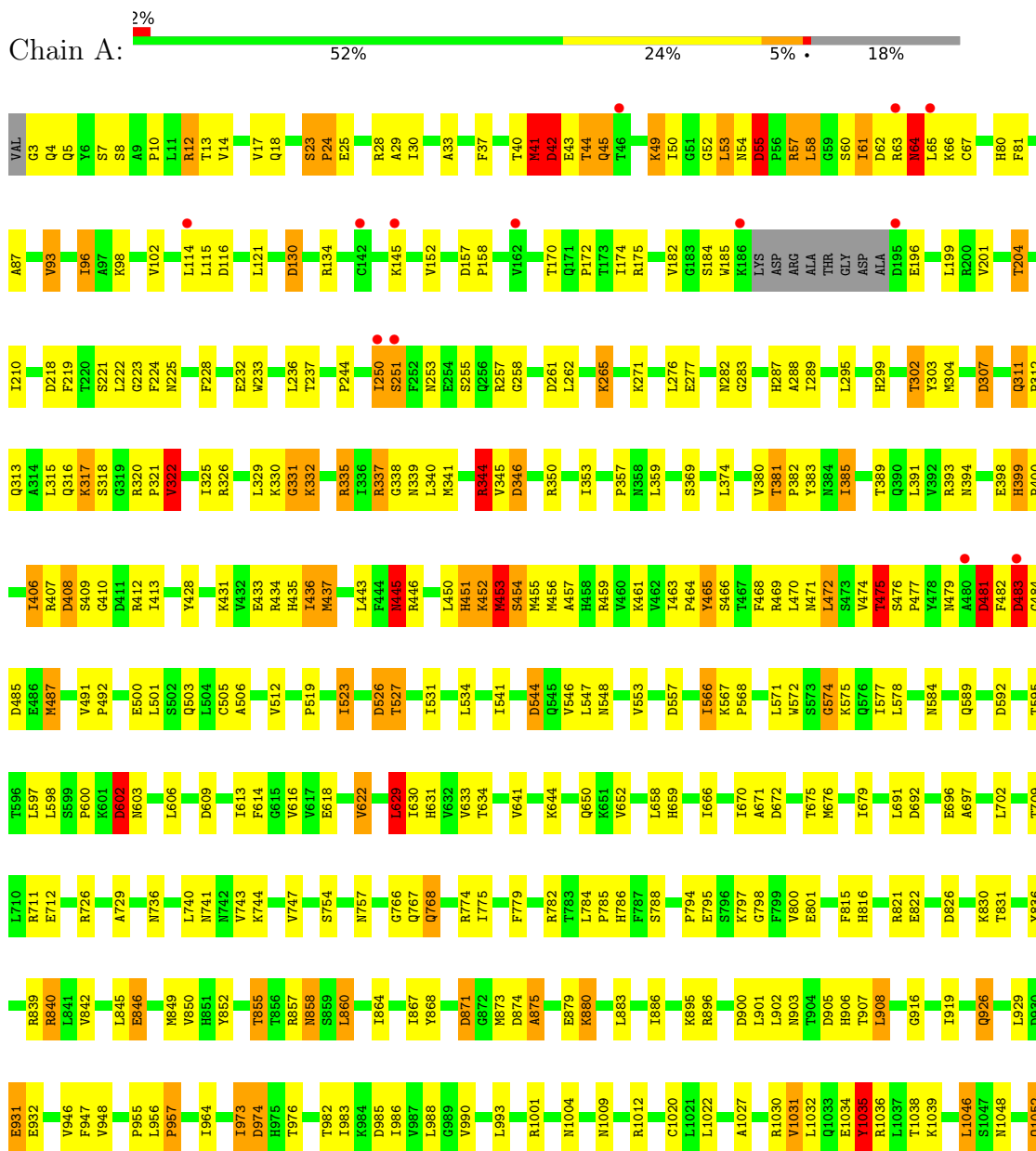
- Molecule 15 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	A	1	Total 1	Mg 1	0	0

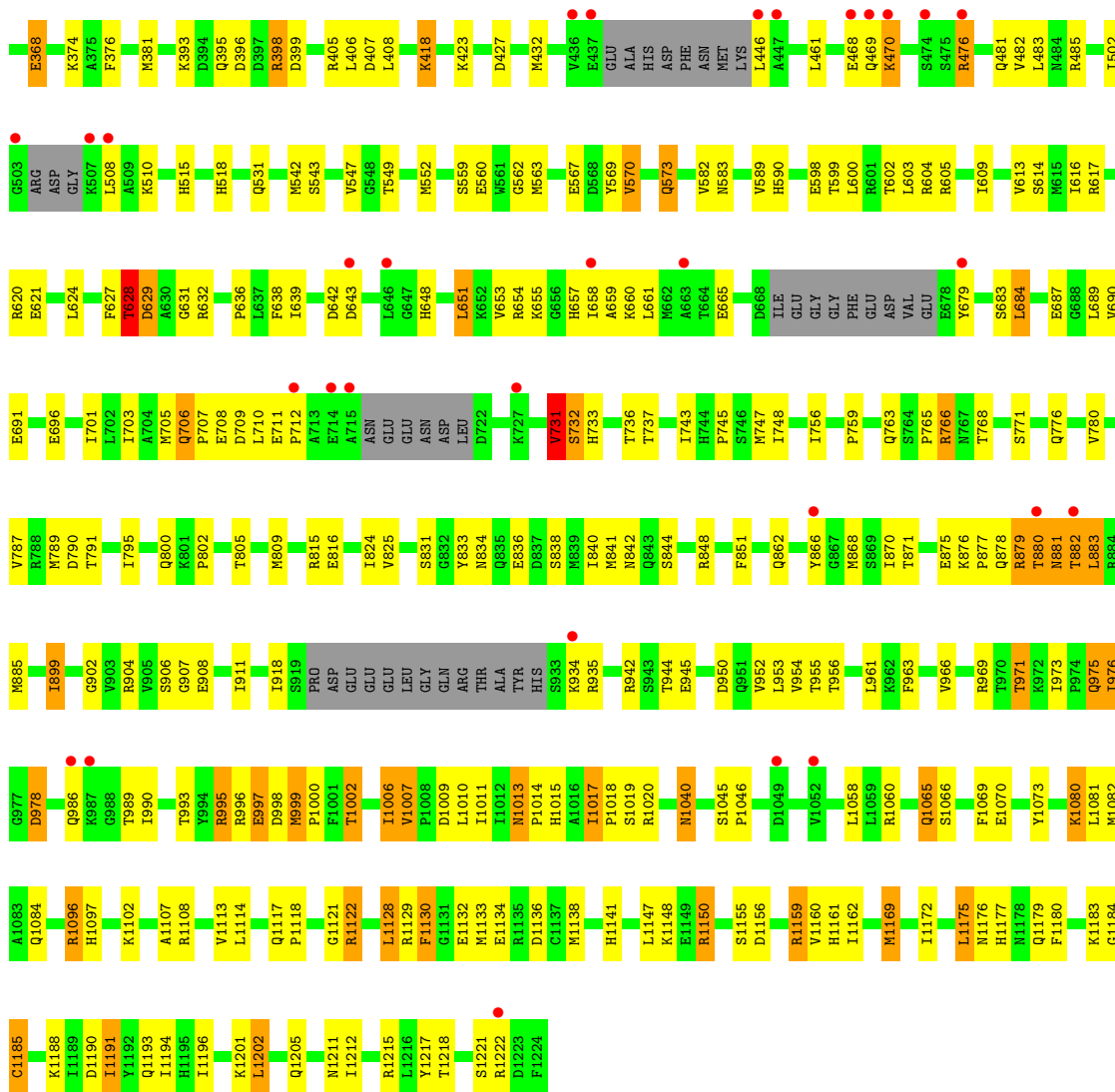
### 3 Residue-property plots

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 electron 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 dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). 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.

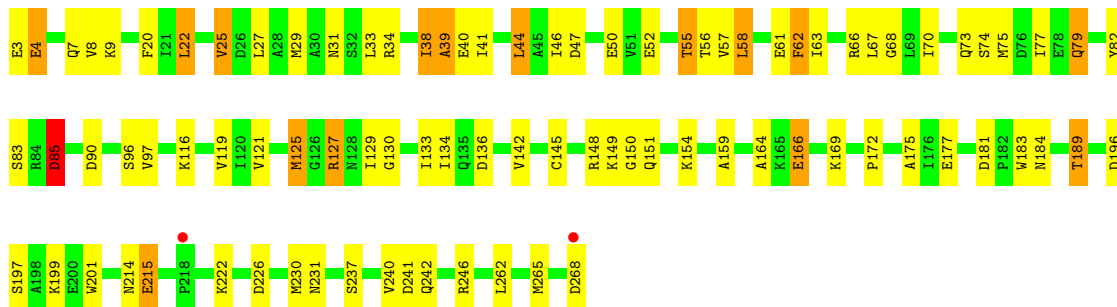
- Molecule 1: DNA-directed RNA polymerase II subunit RPB1





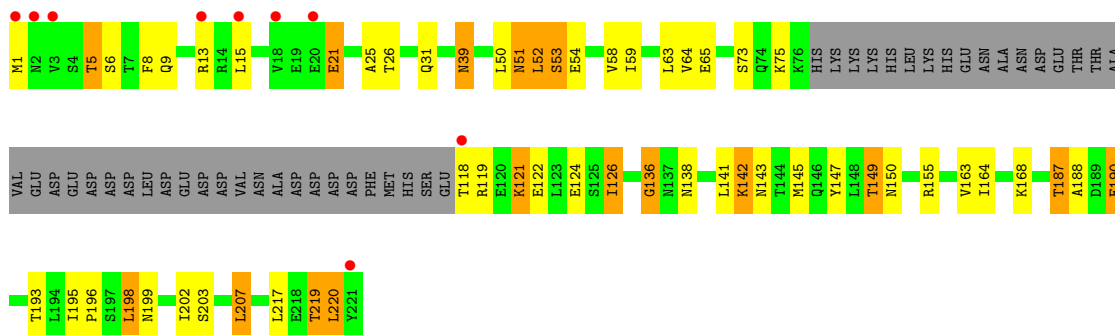


● Molecule 3: DNA-directed RNA polymerase II subunit RPB3

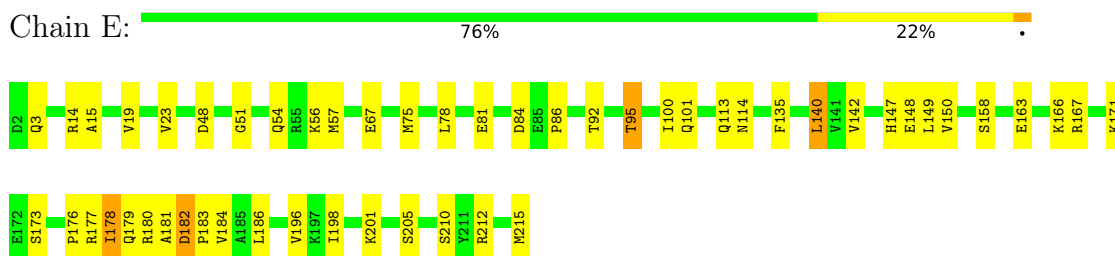


● Molecule 4: DNA-directed RNA polymerase II subunit RPB4

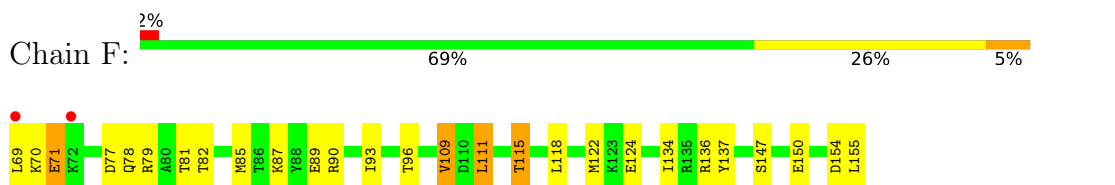




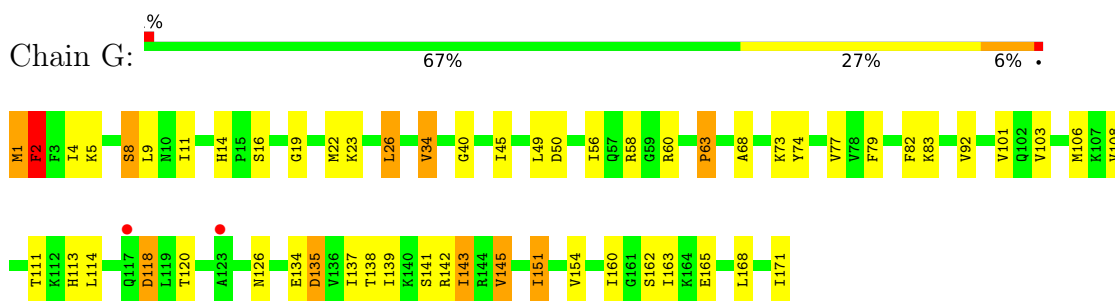
- Molecule 5: DNA-directed RNA polymerases I, II, and III subunit RPABC1



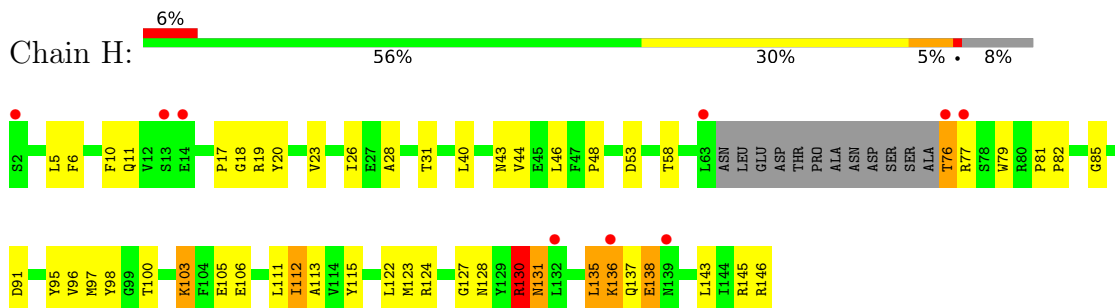
- Molecule 6: DNA-directed RNA polymerases I, II, and III subunit RPABC2



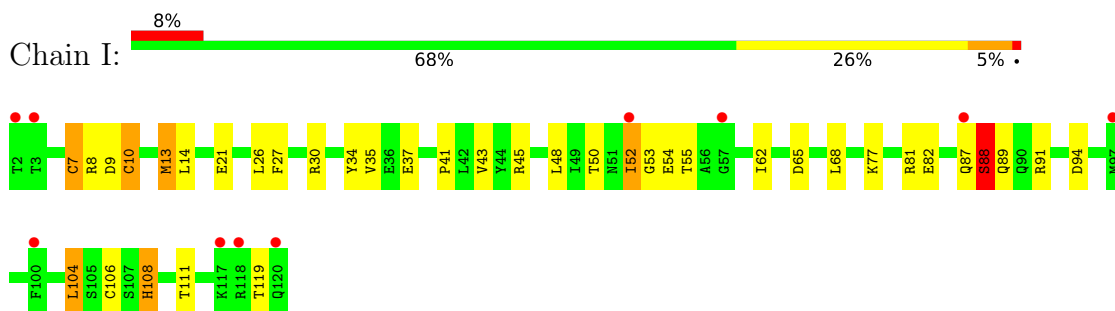
- Molecule 7: DNA-directed RNA polymerase II subunit RPB7



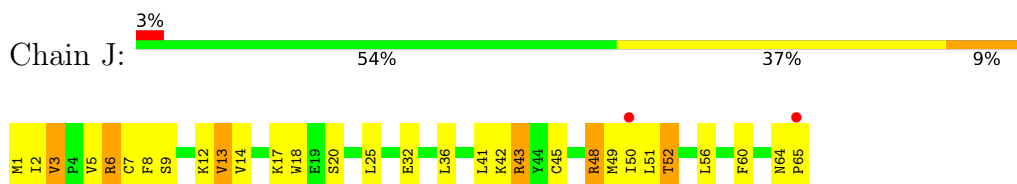
- Molecule 8: DNA-directed RNA polymerases I, II, and III subunit RPABC3



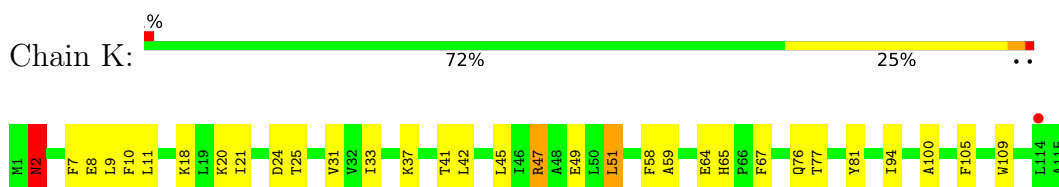
- Molecule 9: DNA-directed RNA polymerase II subunit RPB9



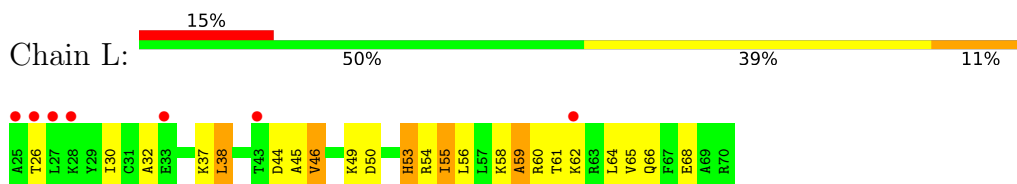
- Molecule 10: DNA-directed RNA polymerases I, II, and III subunit RPABC5



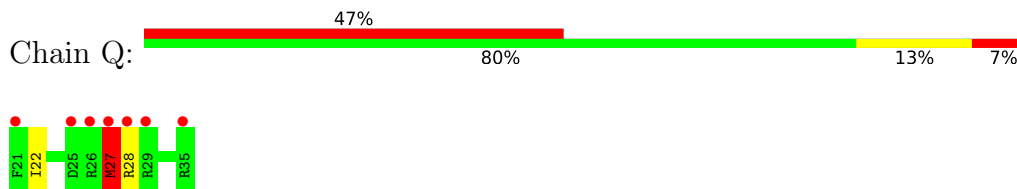
- Molecule 11: DNA-directed RNA polymerase II subunit RPB11



- Molecule 12: DNA-directed RNA polymerases I, II, and III subunit RPABC4



- Molecule 13: PHE-ILE-LYS-ARG-ASP-ARG-MET-ARG-ARG-ASN-PHE-LEU-ARG-MET-ARG



## 4 Data and refinement statistics i

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	222.82Å 392.73Å 283.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.93 – 3.90 48.93 – 3.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.93-3.90) 99.9 (48.93-3.90)	Depositor EDS
$R_{merge}$	0.21	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.33 (at 3.88Å)	Xtrriage
Refinement program	BUSTER-TNT 2.10.2	Depositor
R, $R_{free}$	0.162 , 0.194 0.182 , 0.213	Depositor DCC
$R_{free}$ test set	2265 reflections (2.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	129.8	Xtrriage
Anisotropy	0.419	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 140.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.017 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.024 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	31339	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	130.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.85% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 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

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	0.91	9/11391 (0.1%)	1.50	120/15402 (0.8%)
2	B	0.87	2/9048 (0.0%)	1.43	73/12200 (0.6%)
3	C	0.81	0/2133	1.38	14/2891 (0.5%)
4	D	0.86	0/1450	1.64	33/1945 (1.7%)
5	E	0.78	0/1788	1.38	11/2406 (0.5%)
6	F	0.90	0/717	1.47	8/967 (0.8%)
7	G	0.85	0/1368	1.34	10/1844 (0.5%)
8	H	0.77	0/1086	1.31	10/1470 (0.7%)
9	I	0.84	0/989	1.41	6/1331 (0.5%)
10	J	0.87	0/541	1.46	3/727 (0.4%)
11	K	0.75	0/938	1.39	6/1267 (0.5%)
12	L	0.78	0/365	1.38	3/485 (0.6%)
13	Q	3.63	2/77 (2.6%)	2.16	3/105 (2.9%)
All	All	0.88	13/31891 (0.0%)	1.45	300/43040 (0.7%)

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	Q	27	MET	SD-CE	28.20	2.50	1.79
1	A	57	ARG	CA-C	10.60	1.67	1.52
13	Q	27	MET	CG-SD	8.32	2.01	1.80
1	A	57	ARG	CA-CB	6.50	1.64	1.53
2	B	882	THR	CA-C	6.29	1.60	1.52

The worst 5 of 300 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	483	ASP	CA-CB-CG	15.05	127.65	112.60
13	Q	28	ARG	N-CA-C	-13.03	95.66	111.69
1	A	957	PRO	CA-C-N	-11.48	107.83	123.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	957	PRO	C-N-CA	-11.48	107.83	123.10
1	A	55	ASP	CA-CB-CG	10.64	123.24	112.60

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	11190	0	11255	256	0
2	B	8876	0	8919	165	0
3	C	2095	0	2051	44	0
4	D	1440	0	1456	16	0
5	E	1752	0	1776	22	0
6	F	705	0	731	16	0
7	G	1340	0	1357	25	0
8	H	1068	0	1040	25	0
9	I	971	0	927	17	0
10	J	532	0	542	20	0
11	K	920	0	929	22	0
12	L	363	0	386	10	0
13	Q	78	0	36	3	0
14	A	2	0	0	0	0
14	B	1	0	0	0	0
14	C	1	0	0	0	0
14	I	2	0	0	0	0
14	J	1	0	0	0	0
14	L	1	0	0	0	0
15	A	1	0	0	0	0
All	All	31339	0	31405	556	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 9.

The worst 5 of 556 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)
13:Q:27:MET:CG	13:Q:27:MET:SD	2.01	1.47
1:A:868:TYR:CE1	1:A:1064:VAL:HG11	1.85	1.10
1:A:1081:LEU:HB3	1:A:1082:ASN:HA	1.40	1.04
2:B:121:ASN:HD22	2:B:207:GLY:HA3	1.24	1.02
13:Q:27:MET:SD	13:Q:27:MET:CE	2.50	0.99

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 X-ray entries followed by that with respect to entries of similar resolution.

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	1411/1732 (82%)	1237 (88%)	140 (10%)	34 (2%)	4	29
2	B	1102/1223 (90%)	961 (87%)	114 (10%)	27 (2%)	4	29
3	C	264/266 (99%)	240 (91%)	20 (8%)	4 (2%)	8	37
4	D	176/221 (80%)	156 (89%)	15 (8%)	5 (3%)	4	27
5	E	212/214 (99%)	196 (92%)	14 (7%)	2 (1%)	14	47
6	F	85/87 (98%)	79 (93%)	6 (7%)	0	100	100
7	G	169/171 (99%)	159 (94%)	8 (5%)	2 (1%)	10	41
8	H	129/145 (89%)	110 (85%)	14 (11%)	5 (4%)	2	21
9	I	117/119 (98%)	100 (86%)	15 (13%)	2 (2%)	7	35
10	J	63/65 (97%)	54 (86%)	6 (10%)	3 (5%)	2	18
11	K	113/115 (98%)	108 (96%)	4 (4%)	1 (1%)	14	47
12	L	44/46 (96%)	30 (68%)	10 (23%)	4 (9%)	0	9
13	Q	13/15 (87%)	11 (85%)	2 (15%)	0	100	100
All	All	3898/4419 (88%)	3441 (88%)	368 (9%)	89 (2%)	5	30

5 of 89 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	ASP
1	A	335	ARG
1	A	775	ILE
2	B	67	SER
2	B	339	THR

### 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 X-ray entries followed by that with respect to entries of similar resolution.

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	1244/1519 (82%)	1047 (84%)	197 (16%)	2	15
2	B	967/1060 (91%)	822 (85%)	145 (15%)	3	16
3	C	234/234 (100%)	196 (84%)	38 (16%)	2	14
4	D	160/200 (80%)	134 (84%)	26 (16%)	2	14
5	E	196/196 (100%)	176 (90%)	20 (10%)	7	26
6	F	77/77 (100%)	67 (87%)	10 (13%)	4	19
7	G	152/152 (100%)	131 (86%)	21 (14%)	3	18
8	H	117/127 (92%)	104 (89%)	13 (11%)	6	23
9	I	113/113 (100%)	96 (85%)	17 (15%)	3	16
10	J	60/60 (100%)	51 (85%)	9 (15%)	3	16
11	K	99/99 (100%)	90 (91%)	9 (9%)	9	31
12	L	40/40 (100%)	28 (70%)	12 (30%)	0	2
13	Q	1/15 (7%)	0	1 (100%)	0	0
All	All	3460/3892 (89%)	2942 (85%)	518 (15%)	3	16

5 of 518 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
8	H	76	THR
9	I	43	VAL
8	H	53	ASP
1	A	1445	ILE
1	A	1406	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 93 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	1013	ASN
3	C	184	ASN
2	B	1112	GLN
3	C	54	ASN
4	D	132	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 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

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	1421/1732 (82%)	-0.06	27 (1%) 66 46	55, 112, 198, 278	0
2	B	1118/1223 (91%)	0.11	43 (3%) 44 32	59, 127, 216, 300	0
3	C	266/266 (100%)	-0.12	2 (0%) 82 64	76, 117, 173, 219	0
4	D	180/221 (81%)	0.13	9 (5%) 34 26	92, 131, 203, 241	0
5	E	214/214 (100%)	0.08	0 100 100	86, 162, 233, 254	0
6	F	87/87 (100%)	-0.21	2 (2%) 61 43	59, 89, 131, 142	0
7	G	171/171 (100%)	-0.15	2 (1%) 76 55	79, 111, 158, 224	0
8	H	133/145 (91%)	0.21	9 (6%) 23 21	104, 166, 223, 237	0
9	I	119/119 (100%)	0.38	10 (8%) 17 18	128, 157, 227, 246	0
10	J	65/65 (100%)	0.07	2 (3%) 51 36	85, 111, 162, 184	0
11	K	115/115 (100%)	-0.26	1 (0%) 81 62	66, 113, 165, 200	0
12	L	46/46 (100%)	0.68	7 (15%) 5 8	98, 169, 211, 238	0
13	Q	15/15 (100%)	2.04	7 (46%) 0 1	108, 188, 212, 222	0
All	All	3950/4419 (89%)	0.03	121 (3%) 51 36	55, 122, 210, 300	0

The worst 5 of 121 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1086	PHE	9.3
4	D	3	VAL	6.0
1	A	65	LEU	5.4
2	B	446	LEU	5.1
12	L	27	LEU	5.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
15	MG	A	1803	1/1	0.72	0.38	50,50,50,50	0
14	ZN	L	101	1/1	0.99	0.03	149,149,149,149	0
14	ZN	I	202	1/1	0.99	0.04	211,211,211,211	0
14	ZN	C	301	1/1	1.00	0.03	94,94,94,94	0
14	ZN	I	201	1/1	1.00	0.03	127,127,127,127	0
14	ZN	A	1801	1/1	1.00	0.02	163,163,163,163	0
14	ZN	J	101	1/1	1.00	0.05	111,111,111,111	0
14	ZN	A	1802	1/1	1.00	0.04	97,97,97,97	0
14	ZN	B	1301	1/1	1.00	0.07	112,112,112,112	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.