



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 4BOC  
Title : Structure of mitochondrial RNA polymerase elongation complex  
Authors : Schwinghammer, K.; Cheung, A.; Morozov, Y.; Agaronyan, K.; Temiakov, D.;  
Cramer, P.  
Deposited on : 2013-05-18  
Resolution : 2.65 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.

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

MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

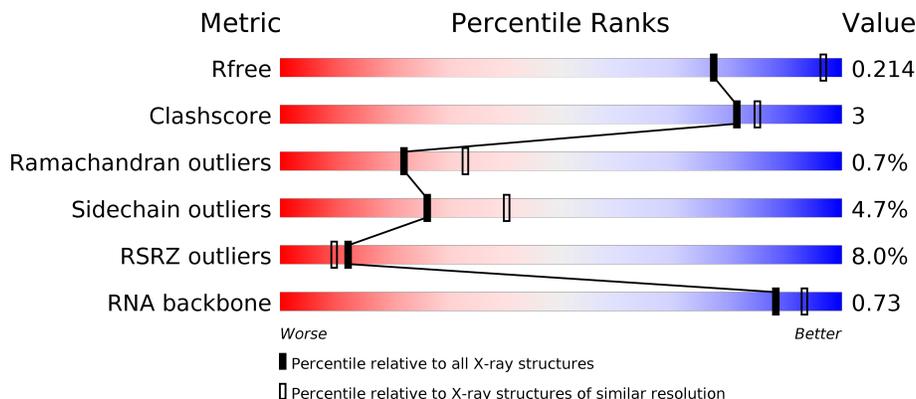
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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))
$R_{free}$	130704	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)
RNA backbone	3102	1010 (2.96-2.36)

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 on 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	1088	
2	N	28	
3	R	14	
4	T	28	

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 9302 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, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	989	7880	5017	1423	1392	48	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	MET	-	expression tag	UNP O00411
A	144	GLY	-	expression tag	UNP O00411
A	145	HIS	-	expression tag	UNP O00411
A	146	HIS	-	expression tag	UNP O00411
A	147	HIS	-	expression tag	UNP O00411
A	148	HIS	-	expression tag	UNP O00411
A	149	HIS	-	expression tag	UNP O00411
A	150	HIS	-	expression tag	UNP O00411

- Molecule 2 is a DNA chain called 5'-D(\*CP\*AP\*TP\*GP\*GP\*GP\*GP\*TP\*AP\*AP\*TP\*T  
P\*AP\*TP \*TP\*TP\*CP\*GP\*AP\*CP\*GP\*CP\*CP\*AP\*GP\*AP\*CP\*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	N	21	437	205	89	122	21	0	0	0

- Molecule 3 is a RNA chain called 5'-R(\*AP\*GP\*UP\*CP\*UP\*GP\*CP\*GP\*GP\*CP\*GP\*C  
P\*GP\*CP)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	R	9	195	86	37	63	9	0	0	0

- Molecule 4 is a DNA chain called 5'-D(\*CP\*GP\*TP\*CP\*TP\*GP\*GP\*CP\*GP\*TP\*GP\*C  
P\*GP\*CP \*GP\*CP\*CP\*GP\*CP\*TP\*AP\*CP\*CP\*CP\*CP\*AP\*TP\*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
4	T	27	546	258	96	165	27	0	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	227	Total 227	O 227	0	0
5	N	7	Total 7	O 7	0	0
5	R	5	Total 5	O 5	0	0
5	T	5	Total 5	O 5	0	0





- Molecule 4: 5'-D(\*CP\*GP\*TP\*CP\*TP\*GP\*GP\*CP\*GP\*TP\*GP\*CP\*GP\*CP \*GP\*CP\*CP\*GP\*CP\*TP\*AP\*CP\*CP\*CP\*CP\*AP\*TP\*G)-3'



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	225.19Å 225.19Å 225.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.81 – 2.65 39.81 – 2.65	Depositor EDS
% Data completeness (in resolution range)	100.0 (39.81-2.65) 100.0 (39.81-2.65)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.83 (at 2.65Å)	Xtrriage
Refinement program	BUSTER 2.11.5	Depositor
R, $R_{free}$	0.179 , 0.209 0.186 , 0.214	Depositor DCC
$R_{free}$ test set	2755 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	74.8	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 68.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.023 for -l,-k,-h	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9302	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	97.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.23% 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](#)

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	0.44	0/8071	0.65	0/10952
2	N	1.07	0/491	0.94	0/754
3	R	1.13	0/217	0.96	0/337
4	T	1.07	0/609	1.00	3/936 (0.3%)
All	All	0.58	0/9388	0.71	3/12979 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	T	0	DT	O4'-C4'-C3'	-6.48	101.91	104.50
4	T	4	DC	C4'-C3'-C2'	-5.12	98.49	103.10
4	T	0	DT	P-O3'-C3'	5.09	125.81	119.70

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	7880	0	7966	40	0
2	N	437	0	235	2	0
3	R	195	0	100	2	0
4	T	546	0	303	7	0
5	A	227	0	0	1	0
5	N	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	R	5	0	0	0	0
5	T	5	0	0	0	0
All	All	9302	0	8604	45	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 3.

The worst 5 of 45 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)
1:A:419:LYS:HB2	1:A:420:PRO:HA	1.39	1.04
1:A:363:GLN:HB3	1:A:364:LEU:HG	1.60	0.83
1:A:419:LYS:CB	1:A:420:PRO:HA	2.13	0.78
1:A:747:PRO:HB2	1:A:748:HIS:HA	1.71	0.71
1:A:419:LYS:HB2	1:A:420:PRO:CA	2.19	0.68

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	983/1088 (90%)	947 (96%)	29 (3%)	7 (1%)	22 33

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	419	LYS
1	A	600	SER
1	A	615	VAL
1	A	662	ALA
1	A	673	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	857/943 (91%)	817 (95%)	40 (5%)	26 40

5 of 40 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	668	THR
1	A	683	LEU
1	A	1197	LYS
1	A	679	GLN
1	A	758	LEU

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

Mol	Chain	Res	Type
1	A	331	GLN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	R	9/14 (64%)	1 (11%)	1 (11%)

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	R	2	C

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	R	1	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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 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	989/1088 (90%)	0.33	76 (7%) 13   10	51, 84, 152, 198	0
2	N	21/28 (75%)	0.37	6 (28%) 0   0	87, 145, 199, 230	0
3	R	9/14 (64%)	-0.21	0 100   100	63, 72, 105, 116	0
4	T	27/28 (96%)	0.24	2 (7%) 14   12	73, 111, 233, 238	0
All	All	1046/1158 (90%)	0.32	84 (8%) 12   9	51, 85, 158, 238	0

The worst 5 of 84 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	218	GLN	7.7
1	A	745	HIS	7.2
1	A	749	SER	6.8
1	A	744	ALA	6.7
1	A	364	LEU	6.3

### 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 carbohydrates in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.