



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 01:18 AM GMT

PDB ID : 2CKZ  
Title : X-RAY STRUCTURE OF RNA POLYMERASE III SUBCOMPLEX C17-C25.  
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Deposited on : 2006-04-24  
Resolution : 3.20 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  
<http://wwpdb.org/validation/2016/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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

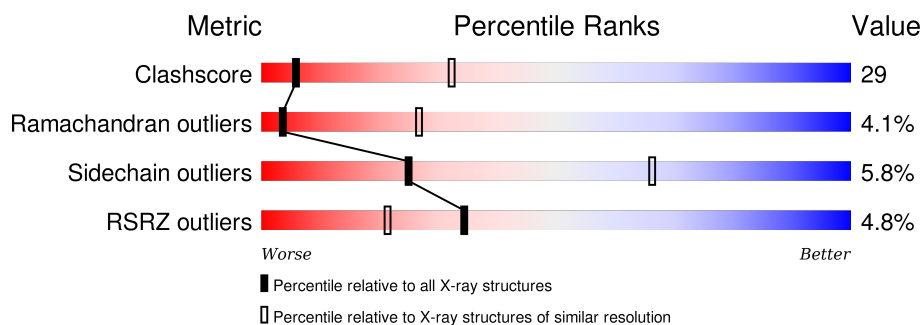
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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))
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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. 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	161	<div> <div>3%</div> <div>16% 19% .. 63%</div> </div>
1	C	161	<div> <div>4%</div> <div>42% 32% • 22%</div> </div>
2	B	218	<div> <div>%</div> <div>45% 33% • 17%</div> </div>
2	D	218	<div> <div>6%</div> <div>38% 38% 5% • 19%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4386 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.

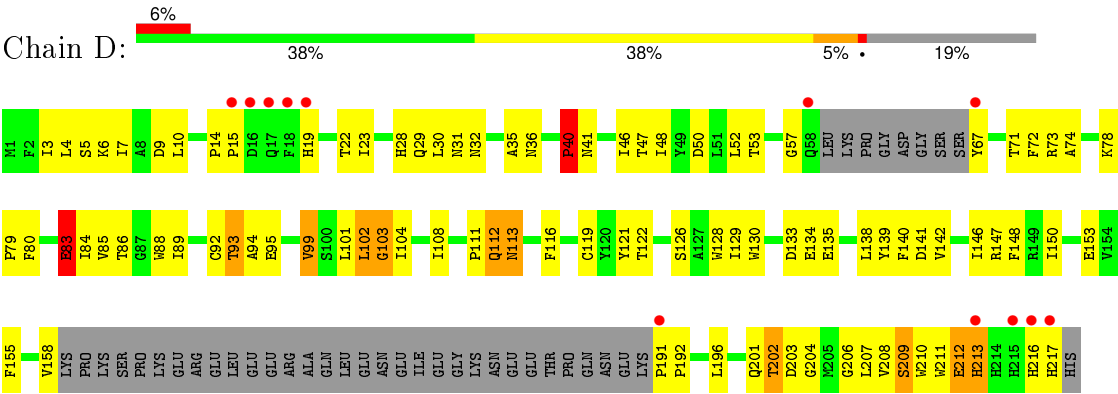
- Molecule 1 is a protein called DNA-DIRECTED RNA POLYMERASE III 18 KD POLYPEPTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	59	Total	C	N	O	S	0	0	1
			487	314	84	88	1			
1	C	125	Total	C	N	O	S	0	0	1
			1022	657	166	193	6			

- Molecule 2 is a protein called DNA-DIRECTED RNA POLYMERASE III 25 KD POLYPEPTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	180	Total	C	N	O	S	0	0	1
			1440	943	229	260	8			
2	D	177	Total	C	N	O	S	0	0	1
			1437	940	232	257	8			





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	138.23Å 138.23Å 247.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.20 29.93 – 3.10	Depositor EDS
% Data completeness (in resolution range)	88.9 (20.00-3.20) 87.1 (29.93-3.10)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.57 (at 3.11Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.236 , 0.307 0.229 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	104.3	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 103.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 22673 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4386	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.61% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.41	0/495	0.57	0/665
1	C	0.39	0/1037	0.56	0/1391
2	B	0.42	0/1477	0.67	0/2006
2	D	0.40	0/1477	0.63	0/2007
All	All	0.41	0/4486	0.62	0/6069

There are no bond length outliers.

There are no bond angle outliers.

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	487	0	494	36	0
1	C	1022	0	1032	56	0
2	B	1440	0	1434	82	0
2	D	1437	0	1411	108	0
All	All	4386	0	4371	252	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 29.

The worst 5 of 252 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)
2:B:93:THR:HG23	2:B:95:GLU:H	1.23	1.02
1:A:20:LEU:HD22	1:A:59:THR:HG23	1.48	0.94
1:C:96:SER:HB3	2:D:203:ASP:O	1.67	0.93
2:B:147:ARG:HH12	2:B:208:VAL:HG22	1.34	0.92
1:C:48:ARG:HA	1:C:49:PRO:N	1.88	0.89

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	54/161 (34%)	42 (78%)	10 (18%)	2 (4%)	4	29
1	C	118/161 (73%)	94 (80%)	20 (17%)	4 (3%)	5	31
2	B	174/218 (80%)	153 (88%)	17 (10%)	4 (2%)	8	44
2	D	171/218 (78%)	142 (83%)	18 (10%)	11 (6%)	2	13
All	All	517/758 (68%)	431 (83%)	65 (13%)	21 (4%)	3	27

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	19	HIS
2	D	40	PRO
2	D	102	LEU
2	D	213	HIS
2	B	64	GLY

### 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	54/145 (37%)	52 (96%)	2 (4%)	41	79
1	C	115/145 (79%)	110 (96%)	5 (4%)	35	75
2	B	159/196 (81%)	148 (93%)	11 (7%)	19	59
2	D	158/196 (81%)	148 (94%)	10 (6%)	22	63
All	All	486/682 (71%)	458 (94%)	28 (6%)	25	66

5 of 28 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	211	TRP
1	C	51	ASN
2	D	139	TYR
1	C	17	LEU
1	C	21	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	51	ASN
1	C	52	HIS
2	D	112	GLN
1	C	31	GLN
2	D	113	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

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	59/161 (36%)	0.05	5 (8%)	13 7	56, 100, 187, 204	0
1	C	125/161 (77%)	-0.14	6 (4%)	34 21	61, 112, 196, 208	0
2	B	180/218 (82%)	-0.19	3 (1%)	73 60	55, 84, 151, 206	0
2	D	177/218 (81%)	0.05	12 (6%)	20 11	52, 95, 207, 231	0
All	All	541/758 (71%)	-0.07	26 (4%)	34 21	52, 95, 188, 231	0

The worst 5 of 26 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	67	TYR	5.0
1	A	37	LEU	4.8
2	D	216	HIS	4.5
2	D	19	HIS	4.2
2	D	15	PRO	4.0

### 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 carbohydrates in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.