

# wwPDB X-ray Structure Validation Summary Report (i)

#### Oct 31, 2023 – 06:41 PM EDT

PDB ID : 3U28

Title : Crystal structure of a Cbf5-Nop10-Gar1 complex from Saccharomyces cere-

visiae

 $\begin{array}{cccc} \text{Authors} & : & \text{Ye, K.; Li, S.} \\ \text{Deposited on} & : & 2011\text{-}10\text{-}02 \end{array}$ 

Resolution : 1.90 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*A user guide is available at

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

MolProbity: 4.02b-467 Xtriage (Phenix): 1.13

EDS: 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

Validation Pipeline (wwPDB-VP) : 2.36

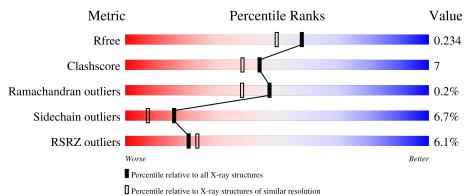
## 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$ 

The reported resolution of this entry is 1.90 Å.

RSRZ outliers

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.



127900

Whole archive Similar resolution Metric (#Entries) (#Entries, resolution range(Å)) $R_{free}$ 6207 (1.90-1.90) 130704 Clashscore 141614 6847 (1.90-1.90) Ramachandran outliers 138981 6760 (1.90-1.90) Sidechain outliers 138945 6760 (1.90-1.90)

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 >=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 <=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.

6082 (1.90-1.90)

Mol	Chain	Length	Quality of chain							
1	A	400	70%	13%		16%				
2	В	58	79%			17%				
3	С	114	68%	11% •		19%				



## 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4028 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 H/ACA ribonucleoprotein complex subunit 4.

$\mathbf{Mol}$	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	A	335	Total 2641	C 1667	N 470	O 487	S 17	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP P33322
A	2	ALA	-	EXPRESSION TAG	UNP P33322
A	395	HIS	-	EXPRESSION TAG	UNP P33322
A	396	HIS	-	EXPRESSION TAG	UNP P33322
A	397	HIS	-	EXPRESSION TAG	UNP P33322
A	398	HIS	-	EXPRESSION TAG	UNP P33322
A	399	HIS	-	EXPRESSION TAG	UNP P33322
A	400	HIS	-	EXPRESSION TAG	UNP P33322

• Molecule 2 is a protein called H/ACA ribonucleoprotein complex subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	В	48	Total 387	C 244	N 70	O 71	S	0	0	0

• Molecule 3 is a protein called H/ACA ribonucleoprotein complex subunit 1.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
3	С	92	Total 721	C 470	N 113	O 134	S 4	0	0	0

There are 21 discrepancies between the modelled and reference sequences:

Cl	hain	Residue	Modelled	Actual	Comment	Reference
	С	11	MET	-	EXPRESSION TAG	UNP P28007
	С	12	GLY	-	EXPRESSION TAG	UNP P28007

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Chain	Residue	Modelled	Actual	Comment	Reference
С	13	SER	-	EXPRESSION TAG	UNP P28007
С	14	SER	-	EXPRESSION TAG	UNP P28007
С	15	HIS	-	EXPRESSION TAG	UNP P28007
С	16	HIS	-	EXPRESSION TAG	UNP P28007
С	17	HIS	-	EXPRESSION TAG	UNP P28007
С	18	HIS	-	EXPRESSION TAG	UNP P28007
С	19	HIS	-	EXPRESSION TAG	UNP P28007
С	20	HIS	-	EXPRESSION TAG	UNP P28007
С	21	SER	-	EXPRESSION TAG	UNP P28007
С	22	SER	-	EXPRESSION TAG	UNP P28007
С	23	GLY	-	EXPRESSION TAG	UNP P28007
С	24	LEU	-	EXPRESSION TAG	UNP P28007
С	25	VAL	-	EXPRESSION TAG	UNP P28007
С	26	PRO	-	EXPRESSION TAG	UNP P28007
С	27	ARG	-	EXPRESSION TAG	UNP P28007
С	28	GLY	-	EXPRESSION TAG	UNP P28007
С	29	SER	-	EXPRESSION TAG	UNP P28007
С	30	HIS	-	EXPRESSION TAG	UNP P28007
С	31	MET	-	EXPRESSION TAG	UNP P28007

#### • Molecule 4 is water.

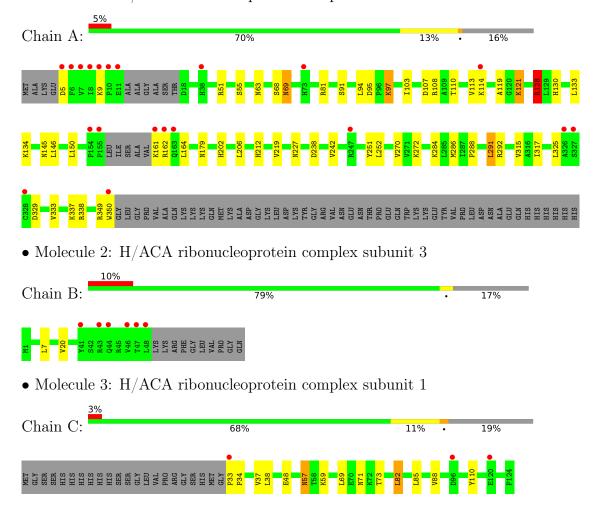
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	183	Total O 183 183	0	0
4	В	44	Total O 44 44	0	0
4	С	52	Total O 52 52	0	0



## 3 Residue-property plots (i)

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: H/ACA ribonucleoprotein complex subunit 4





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	54.41Å 59.39Å 169.21Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 - 1.90	Depositor
Resolution (A)	19.92 - 1.90	EDS
% Data completeness	89.7 (19.92-1.90)	Depositor
(in resolution range)	89.7 (19.92-1.90)	EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	4.13 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
D D.	0.204 , 0.233	Depositor
$R, R_{free}$	0.204 , $0.234$	DCC
$R_{free}$ test set	1971 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.1	Xtriage
Anisotropy	0.367	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.37, 43.9	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4028	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 8.29% of the height of the origin peak. No significant pseudotranslation is detected.

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



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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		
MIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.50	0/2686	0.66	3/3630 (0.1%)	
2	В	0.43	0/395	0.59	0/530	
3	С	0.54	0/739	0.68	0/1002	
All	All	0.50	0/3820	0.65	3/5162 (0.1%)	

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
1	A	128	ARG	NE-CZ-NH1	6.57	123.59	120.30
1	A	108	ARG	NE-CZ-NH2	-6.03	117.29	120.30
1	A	128	ARG	NE-CZ-NH2	-5.70	117.45	120.30

There are no chirality outliers.

There are no planarity outliers.

#### 5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2641	0	2695	43	0
2	В	387	0	398	0	0
3	С	721	0	721	10	0
4	A	183	0	0	8	0
4	В	44	0	0	0	0
4	С	52	0	0	2	0
All	All	4028	0	3814	50	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.

The worst 5 of 50 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:133:LEU:H	1:A:179:ASN:HD21	1.20	0.89
3:C:57:ASN:HD22	3:C:59:LYS:H	1.21	0.87
1:A:270:VAL:CG1	1:A:292:ARG:HB2	2.08	0.84
1:A:130:HIS:HE1	1:A:212:HIS:ND1	1.77	0.81
3:C:37:VAL:HG22	3:C:110:TYR:HB3	1.64	0.80

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	Percentil	les
1	A	329/400 (82%)	320 (97%)	8 (2%)	1 (0%)	41 31	
2	В	46/58 (79%)	46 (100%)	0	0	100 10	)()
3	С	90/114 (79%)	85 (94%)	5 (6%)	0	100 10	00
All	All	465/572 (81%)	451 (97%)	13 (3%)	1 (0%)	47 38	3

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	69	ASN

#### 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	293/345~(85%)	273 (93%)	20 (7%)	16 7
2	В	43/51 (84%)	41 (95%)	2 (5%)	26 16
3	С	80/98 (82%)	74 (92%)	6 (8%)	13 5
All	All	416/494 (84%)	388 (93%)	28 (7%)	16 7

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

Mol	Chain	Res	Type
1	A	164	LEU
3	С	88	VAL
1	A	252	LEU
3	С	69	LEU
1	A	251	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
3	С	57	ASN
1	A	245	ASN
1	A	214	GLN
1	A	182	ASN
1	A	234	HIS

#### 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 monosaccharides 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,  $95^{th}$  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	$OWAB(A^2)$	Q < 0.9
1	A	335/400~(83%)	0.23	20 (5%) 21 24	11, 22, 43, 165	0
2	В	48/58 (82%)	0.33	6 (12%) 3 4	11, 26, 71, 80	0
3	С	92/114 (80%)	0.04	3 (3%) 46 49	12, 22, 39, 45	0
All	All	475/572 (83%)	0.21	29 (6%) 21 24	11, 22, 48, 165	0

The worst 5 of 29 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	7	VAL	9.4
1	A	8	ILE	9.2
1	A	6	PHE	9.1
1	A	161	LYS	5.7
1	A	162	ARG	4.8

### 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 monosaccharides 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.

