



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:48 PM GMT

PDB ID : 4WC5  
Title : Structure of tRNA-processing enzyme complex 3  
Authors : Yamashita, S.; Tomita, K.  
Deposited on : 2014-09-04  
Resolution : 3.41 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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

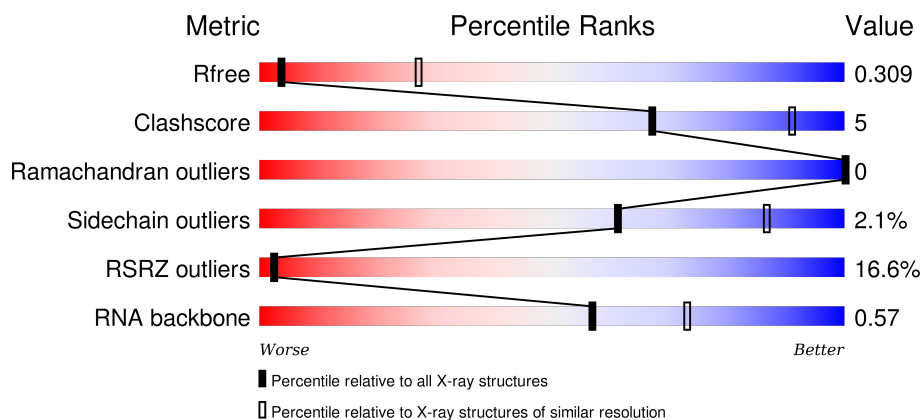
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.41 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))
$R_{free}$	91344	1049 (3.52-3.32)
Clashscore	102246	1032 (3.50-3.34)
Ramachandran outliers	100387	1002 (3.50-3.34)
Sidechain outliers	100360	1003 (3.50-3.34)
RSRZ outliers	91569	1054 (3.52-3.32)
RNA backbone	2183	1042 (4.02-2.80)

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	396	
2	B	74	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4656 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 Poly A polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	0	0	0
			3046	1978	527	534	7			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP O66728
A	384	LYS	-	expression tag	UNP O66728
A	385	LEU	-	expression tag	UNP O66728
A	386	ALA	-	expression tag	UNP O66728
A	387	ALA	-	expression tag	UNP O66728
A	388	ALA	-	expression tag	UNP O66728
A	389	LEU	-	expression tag	UNP O66728
A	390	GLU	-	expression tag	UNP O66728
A	391	HIS	-	expression tag	UNP O66728
A	392	HIS	-	expression tag	UNP O66728
A	393	HIS	-	expression tag	UNP O66728
A	394	HIS	-	expression tag	UNP O66728
A	395	HIS	-	expression tag	UNP O66728
A	396	HIS	-	expression tag	UNP O66728

- Molecule 2 is a RNA chain called RNA (74-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	74	Total	C	N	O	P	0	0	0
			1581	704	283	520	74			

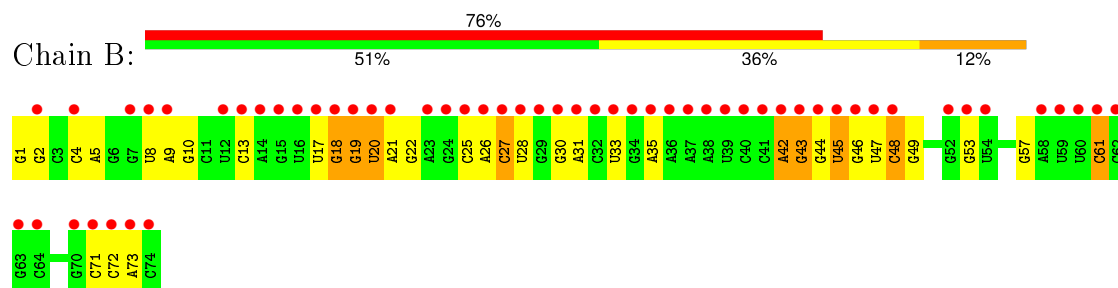
- Molecule 3 is CYTIDINE-5'-TRIPHOSPHATE (three-letter code: CTP) (formula:  $C_9H_{16}N_3O_{14}P_3$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			29	9	3	14	3		



- Molecule 1: Poly A polymerase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.87Å 111.27Å 127.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.82 – 3.41 19.82 – 3.41	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.82-3.41) 99.7 (19.82-3.41)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.40 (at 3.44Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.267 , 0.309 0.267 , 0.309	Depositor DCC
$R_{free}$ test set	592 reflections (4.97%)	DCC
Wilson B-factor (Å <sup>2</sup> )	111.8	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 134.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	1 of 11953 reflections (0.008%)	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4656	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	247.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CTP

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	0.21	0/3097	0.36	0/4150
2	B	0.28	1/1766 (0.1%)	0.67	0/2750
All	All	0.24	1/4863 (0.0%)	0.51	0/6900

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	G	OP3-P	-10.62	1.48	1.61

There are no bond angle outliers.

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	3046	0	3205	28	0
2	B	1581	0	800	14	0
3	A	29	0	12	6	0
All	All	4656	0	4017	41	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 5.

All (41) 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:104:ARG:HA	3:A:501:CTP:O2	1.50	1.08
1:A:104:ARG:CA	3:A:501:CTP:O2	2.30	0.80
2:B:27:C:H42	2:B:43:G:H1	1.38	0.69
1:A:184:LEU:HD23	1:A:220:GLU:HB2	1.81	0.62
1:A:161:GLY:HA3	1:A:208:ILE:HG12	1.83	0.61
1:A:149:ASP:OD2	3:A:501:CTP:N4	2.34	0.60
1:A:372:ILE:HA	1:A:375:LEU:HD12	1.86	0.57
2:B:18:G:O2'	2:B:57:G:N2	2.39	0.55
2:B:42:A:H5'	2:B:43:G:OP2	2.08	0.54
1:A:152:ARG:HE	3:A:501:CTP:HN41	1.56	0.54
1:A:152:ARG:HE	3:A:501:CTP:N4	2.05	0.54
1:A:354:LYS:O	1:A:358:LEU:HG	2.11	0.51
2:B:9:A:O2'	2:B:10:G:N7	2.44	0.51
1:A:293:MET:HA	1:A:297:LEU:HB2	1.92	0.51
1:A:304:LEU:O	1:A:337:LYS:NZ	2.45	0.50
1:A:349:LYS:HB2	1:A:349:LYS:NZ	2.26	0.50
1:A:219:GLU:OE1	1:A:219:GLU:N	2.44	0.49
2:B:33:U:H2'	2:B:35:A:OP2	2.14	0.48
1:A:235:LEU:HD11	1:A:264:LEU:HD21	1.95	0.48
1:A:21:ARG:NH1	1:A:109:ASN:OD1	2.43	0.47
1:A:18:GLY:HA3	3:A:501:CTP:O1B	2.14	0.47
1:A:358:LEU:HD13	1:A:375:LEU:HD13	1.97	0.47
1:A:11:LEU:HD22	1:A:38:GLY:HA3	1.96	0.46
1:A:293:MET:HG3	1:A:297:LEU:HD22	1.97	0.45
1:A:138:ILE:HB	1:A:167:LEU:HD23	1.99	0.45
2:B:27:C:H2'	2:B:28:U:C6	2.52	0.44
1:A:341:TYR:HD1	1:A:345:LEU:HD22	1.82	0.44
2:B:30:G:H2'	2:B:31:A:H8	1.83	0.44
1:A:140:VAL:HG13	1:A:171:THR:HG23	2.00	0.43
2:B:26:A:N6	2:B:45:U:O2	2.50	0.43
1:A:181:LEU:HB2	1:A:183:LEU:HG	2.00	0.43
2:B:53:G:H1	2:B:61:C:H42	1.66	0.43
2:B:19:G:H4'	2:B:20:U:C5'	2.49	0.43
2:B:4:C:H2'	2:B:5:A:O4'	2.18	0.43
1:A:275:HIS:NE2	1:A:279:GLU:OE2	2.51	0.43
1:A:201:ARG:HH22	2:B:2:G:H4'	1.84	0.42
1:A:341:TYR:HA	1:A:345:LEU:HB3	2.01	0.41
1:A:249:PHE:HE2	1:A:342:LEU:HD22	1.85	0.41
2:B:27:C:N3	2:B:43:G:N2	2.49	0.41
1:A:212:TYR:HB3	1:A:218:LEU:HG	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:8:U:O4'	2:B:48:C:O2'	2.38	0.40

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	361/396 (91%)	345 (96%)	16 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	326/349 (93%)	319 (98%)	7 (2%)	61	87

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	230	LYS
1	A	239	ARG
1	A	249	PHE
1	A	294	LYS
1	A	343	GLU

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Mol	Chain	Res	Type
1	A	344	LYS
1	A	349	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	73/74 (98%)	21 (28%)	2 (2%)

All (21) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	13	C
2	B	17	U
2	B	18	G
2	B	19	G
2	B	20	U
2	B	21	A
2	B	22	G
2	B	25	C
2	B	27	C
2	B	42	A
2	B	43	G
2	B	44	G
2	B	45	U
2	B	46	G
2	B	47	U
2	B	48	C
2	B	49	G
2	B	61	C
2	B	71	C
2	B	72	C
2	B	73	A

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	19	G
2	B	45	U

## 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 ⓘ

1 ligand is modelled in this entry.

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	$\# Z  > 2$
3	CTP	A	501	-	21,30,30	0.68	0	31,47,47	1.52	3 (9%)

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
3	CTP	A	501	-	-	0/18/38/38	0/2/2/2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	CTP	PB-O3A-PA	-4.35	120.52	132.73
3	A	501	CTP	PB-O3B-PG	-3.62	120.52	132.67
3	A	501	CTP	C2-N3-C4	3.34	120.32	115.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	CTP	6	0

## 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

### 6.1 Protein, DNA and RNA chains

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	367/396 (92%)	0.28	17 (4%) 36 32	99, 167, 343, 440	0
2	B	74/74 (100%)	3.53	56 (75%) 0 0	242, 345, 435, 507	0
All	All	441/470 (93%)	0.83	73 (16%) 2 2	99, 189, 378, 507	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	36	A	10.2
2	B	20	U	8.9
2	B	9	A	8.8
1	A	367	GLU	8.1
2	B	29	G	7.9
1	A	365	GLY	7.7
2	B	33	U	7.7
2	B	13	C	7.6
2	B	31	A	7.3
2	B	37	A	7.1
2	B	44	G	7.1
2	B	32	C	7.0
2	B	52	G	6.8
1	A	368	LEU	6.2
2	B	46	G	6.0
2	B	53	G	5.9
2	B	8	U	5.8
2	B	60	U	5.8
1	A	381	ASN	5.5
2	B	35	A	5.3
2	B	72	C	5.2
2	B	19	G	5.0
2	B	34	G	5.0
2	B	17	U	4.9

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Mol	Chain	Res	Type	RSRZ
2	B	48	C	4.9
2	B	30	G	4.8
2	B	45	U	4.7
1	A	349	LYS	4.6
2	B	26	A	4.3
2	B	73	A	4.1
2	B	24	G	3.9
2	B	43	G	3.8
2	B	41	C	3.8
2	B	28	U	3.5
1	A	357	GLU	3.5
2	B	25	C	3.4
1	A	61	PHE	3.4
1	A	156	ALA	3.3
2	B	23	A	3.2
2	B	38	A	3.2
1	A	350	LEU	3.2
2	B	16	U	3.2
2	B	21	A	3.2
2	B	71	C	3.2
2	B	74	C	3.1
2	B	39	U	3.0
2	B	70	G	3.0
2	B	42	A	2.9
2	B	12	U	2.9
2	B	54	U	2.9
2	B	47	U	2.8
2	B	18	G	2.7
2	B	63	G	2.7
2	B	58	A	2.7
2	B	64	C	2.6
2	B	15	G	2.5
2	B	4	C	2.5
2	B	62	C	2.5
1	A	366	LYS	2.5
2	B	61	C	2.5
2	B	40	C	2.4
1	A	333	GLU	2.3
1	A	360	LYS	2.3
2	B	7	G	2.3
1	A	351	PRO	2.2
1	A	352	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	354	LYS	2.2
2	B	59	U	2.2
2	B	2	G	2.2
1	A	336	GLU	2.1
1	A	356	GLU	2.1
2	B	27	C	2.0
2	B	14	A	2.0

## 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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	CTP	A	501	29/29	0.88	0.25	-0.15	173,206,293,303	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.