

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

#### Jan 3, 2023 – 04:02 PM EST

PDB ID	:	$7\mathrm{TFZ}$
Title	:	[T:Ag+/Hg2+:T-(pH8.5-pH11; 5s)] Metal-mediated DNA base pair in
		tense grity triangle grown at pH $8.5$ and soaked in pH $11$ for 5s
Authors	:	Lu, B.; Vecchioni, S.; Seeman, N.C.; Sha, R.; Ohayon, Y.P.
Deposited on	:	2022-01-07
Resolution	:	6.13  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:

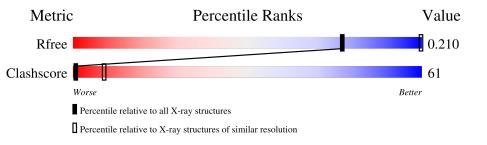
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)		
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

# 1 Overall quality at a glance (i)

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

The reported resolution of this entry is 6.13 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	130704	1003 (8.40-3.88)
Clashscore	141614	1051 (8.30-3.90)

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%

Mol	Chain	Length		Quality of chain					
1	А	21	10%	43%		48%			
2	В	7	14%	14%	71%				
3	С	7	2	9%	57%	14%			
4	D	7		57%		43%			



## 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 858 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 DNA chain called DNA (5'-D(\*GP\*AP\*GP\*CP\*AP\*GP\*CP\*CP\*TP\*GP\* TP\*TP\*GP\*GP\*AP\*CP\*AP\*TP\*CP\*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	А	21	Total 429	C 205	N 80	0 124	Р 20	0	0	0

• Molecule 2 is a DNA chain called DNA (5'-D(P\*CP\*CP\*AP\*TP\*AP\*CP\*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	В	7	Total 140	C 67	N 26	O 40	Р 7	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	С	7	Total 144	C 68	N 25	O 44	P 7	0	0	0
			144	68	ZO	44	1			

• Molecule 4 is a DNA chain called DNA (5'-D(P\*CP\*TP\*GP\*AP\*TP\*GP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	7	Total 145	C 69	N 24	O 45	Р 7	0	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. 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. 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 (5'-D(\*GP\*AP\*GP\*CP\*AP\*GP\*CP\*CP\*TP\*GP\*TP\*TP\*TP\*GP\*GP\*AP \*CP\*AP\*TP\*CP\*A)-3')

Chain A:	10%	43%	48%	, D
61 A2 63 65 65 65 65 65 68 03	19 610 711 712 713 614 615 615 719 719 718 718 718			
• Molecule	2: DNA (5'-D(P	*CP*CP*AP*TP*	AP*CP*A)-3')	
Chain B:	14% 14%		71%	
C1 C2 A3 A5 A7 A7				
• Molecule	e 3: DNA (5'-D(P	*GP*GP*CP*TP*	GP*CP*T)-3')	1
Chain C:	29%	•	57%	14%
68 69 7111 612 714 714				
• Molecule	e 4: DNA (5'-D(P	*CP*TP*GP*AP*	"TP*GP*T)-3')	I
Chain D:	5	7%	43	%
C1 T2 A4 T5 C6 T7				



# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants	107.22Å 107.22Å 85.27Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor
Resolution (Å)	38.74 - 6.13	Depositor
Resolution (A)	53.61 - 5.85	EDS
% Data completeness	70.2 (38.74-6.13)	Depositor
(in resolution range)	58.3(53.61-5.85)	EDS
R <sub>merge</sub>	(Not available)	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	-0.81 (at 5.74Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
P. P.	0.203 , $0.214$	Depositor
R, $R_{free}$	0.203 , $0.210$	DCC
$R_{free}$ test set	34  reflections  (5.44%)	wwPDB-VP
Wilson B-factor $(Å^2)$	156.6	Xtriage
Anisotropy	1.511	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.29, 377.5	EDS
L-test for twinning <sup>2</sup>	$<  L  > = 0.46, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.059 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	858	wwPDB-VP
Average B, all atoms $(Å^2)$	440.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 13.06% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<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	Bo	nd lengths	Bond angles		
	Ullalli	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	1.83	14/481~(2.9%)	1.69	16/741~(2.2%)	
2	В	1.93	3/156~(1.9%)	1.72	4/237~(1.7%)	
3	С	1.14	0/160	1.29	1/245~(0.4%)	
4	D	1.73	3/161~(1.9%)	1.60	5/245~(2.0%)	
All	All	1.74	20/958~(2.1%)	1.62	26/1468~(1.8%)	

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
2	В	3	DA	N9-C4	-11.61	1.30	1.37
4	D	1	DC	OP3-P	-10.74	1.48	1.61
1	А	19	DT	C1'-N1	8.40	1.60	1.49
1	А	11	DT	C3'-O3'	7.90	1.54	1.44
1	А	19	DT	C3'-O3'	7.75	1.54	1.44

The worst 5 of 26 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	4	DC	O4'-C4'-C3'	-8.75	100.75	106.00
4	D	7	DT	O4'-C4'-C3'	-8.13	101.12	106.00
1	А	12	DT	O5'-P-OP1	-8.00	98.50	105.70
1	А	12	DT	O4'-C4'-C3'	-7.91	101.25	106.00
2	В	7	DA	O4'-C1'-N9	7.89	113.52	108.00

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



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	429	0	238	47	0
2	В	140	0	79	25	0
3	С	144	0	80	16	0
4	D	145	0	81	12	0
All	All	858	0	478	78	0

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.

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 61.

The worst 5 of 78 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)
4:D:6:DG:H2"	4:D:7:DT:H5"	1.30	1.07
1:A:13:DT:N3	2:B:3:DA:N1	2.18	0.89
1:A:12:DT:O4	2:B:4:DT:N3	2.16	0.79
2:B:7:DA:H1'	3:C:8:DG:H1'	1.65	0.78
1:A:4:DC:H2'	1:A:5:DA:C8	2.22	0.75

There are no symmetry-related clashes.

### 5.3 Torsion angles (i)

#### 5.3.1 Protein backbone (i)

There are no protein molecules in this entry.

#### 5.3.2 Protein sidechains (i)

There are no protein molecules in this entry.

#### 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)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

#### 6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

