

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

#### Oct 9, 2023 – 11:32 PM EDT

PDB ID : 6XO7

Title: Self-assembly of a 3D DNA crystal lattice (4x6 junction version) containing

the J30 immobile Holliday junction

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Deposited on : 2020-07-06

Resolution : 3.10 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 : 2.35.1

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

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

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

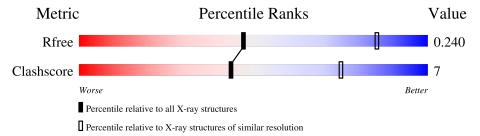
Validation Pipeline (wwPDB-VP) : 2.35.1

# 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 3.10 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}(\mathring{\rm A})) \end{array}$
$R_{free}$	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)

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	A	12	50%	50%					
2	В	9	44%	56%					
3	С	6		100%					
4	D	15	80%		13%	7%			



# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 1333 atoms, of which 474 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\*AP\*CP\*CP\*TP\*GP\*A)-3').

Mol	Chain	Residues		1	Atom	S			ZeroOcc	AltConf	Trace
1	A	12	Total 380	C 117	H 134	N 51	O 67	P 11	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
9	В	0	Total	С	Н	N	О	Р	0	0	0
	Ъ	9	286	87	101	36	53	9	0	U	U

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

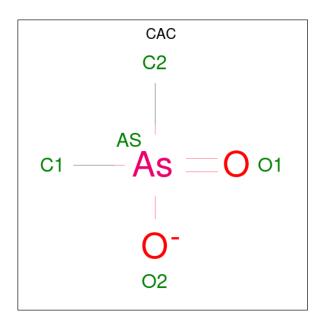
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace			
3	C	6	Total	С	Н	N	О	Р	0	0	0
		U	188	57	68	21	36	6	O	U	O

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

Mol	Chain	Residues		Atoms					ZeroOcc	AltConf	Trace
1	D	15	Total	С	Н	N	О	Р	0	0	0
4	D	10	475	146	171	52	92	14	0	0	

• Molecule 5 is CACODYLATE ION (three-letter code: CAC) (formula: C<sub>2</sub>H<sub>6</sub>AsO<sub>2</sub>).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	В	1	Total As 1 1	0	0
5	D	1	Total As 1 1	0	0

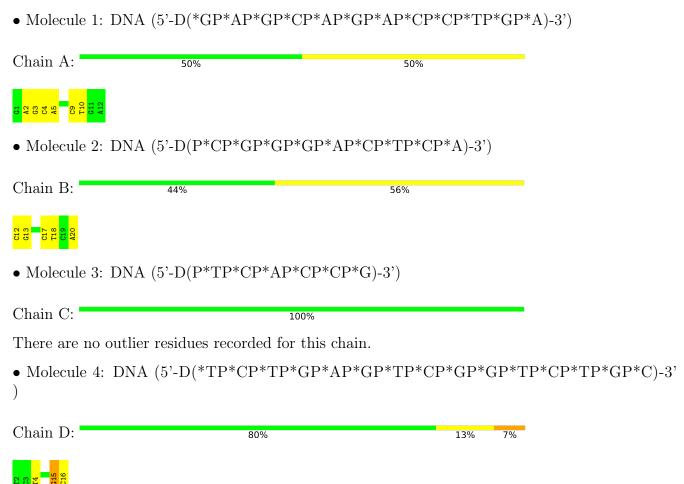
• Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	2	Total Mg 2 2	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.





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32	Depositor
Cell constants	68.14Å 68.14Å 52.77Å	D
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor
Resolution (Å)	34.07 - 3.10	Depositor
Resolution (A)	34.07 - 3.10	EDS
% Data completeness	63.5 (34.07-3.10)	Depositor
(in resolution range)	63.4 (34.07-3.10)	EDS
$R_{merge}$	0.13	Depositor
$R_{sum}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.66 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
$R, R_{free}$	0.215 , $0.240$	Depositor
$\Pi,\ \Pi free$	0.215 , $0.240$	DCC
$R_{free}$ test set	166 reflections $(5.30\%)$	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	62.2	Xtriage
Anisotropy	1.838	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning <sup>2</sup>	$< L > = 0.47, < L^2> = 0.29$	Xtriage
	0.087 for -h,-k,l	
Estimated twinning fraction	0.099  for h,-h-k,-l	Xtriage
	0.105  for -k,-h,-l	
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	1333	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 8.65% 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)

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

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).

Mal	Chain	Bond	lengths	Bond angles		
Mol	Chain	RMSZ $\# Z  > 5$		RMSZ	# Z  > 5	
1	A	0.57	0/277	0.84	0/426	
2	В	0.53	0/207	0.87	0/317	
3	С	0.59	0/133	0.90	0/202	
4	D	0.56	0/339	1.04	1/522~(0.2%)	
All	All	0.56	0/956	0.93	1/1467 (0.1%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
4	D	15	DG	O4'-C1'-N9	5.23	111.66	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 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	246	134	135	3	0
2	В	185	101	101	5	0
3	С	120	68	68	0	0
4	D	304	171	172	3	0
5	В	1	0	0	0	0
5	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	D	2	0	0	0	0
All	All	859	474	476	9	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 9 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
2:B:12:DC:H2'	2:B:13:DG:C8	2.48	0.48
2:B:20:DA:C2	4:D:4:DT:H73	2.48	0.48
2:B:17:DC:H2"	2:B:18:DT:O5'	2.13	0.47
2:B:20:DA:H2	4:D:4:DT:H73	1.78	0.47
2:B:17:DC:H2'	2:B:18:DT:C6	2.51	0.45

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)

Of 4 ligands modelled in this entry, 2 are modelled with single atom and 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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.

