



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 30, 2022 – 04:48 PM EDT

PDB ID : 7JPA
Title : Self-assembly of a 3D DNA crystal lattice (4x6 duplex version) containing the J5 immobile Holliday junction
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Deposited on : 2020-08-07
Resolution : 3.16 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.29
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)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.29

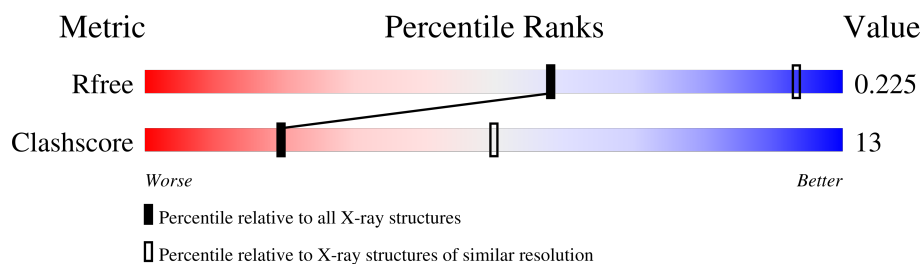
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.16 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}	130704	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	21	<div> <div style="width: 62%; background-color: green;"></div> <div style="width: 38%; background-color: yellow;"></div> </div> <div>62% 38%</div>
2	B	6	<div> <div style="width: 33%; background-color: green;"></div> <div style="width: 67%; background-color: yellow;"></div> </div> <div>33% 67%</div>
3	C	8	<div> <div style="width: 62%; background-color: green;"></div> <div style="width: 38%; background-color: yellow;"></div> </div> <div>62% 38%</div>
4	D	7	<div> <div style="width: 29%; background-color: green;"></div> <div style="width: 71%; background-color: yellow;"></div> </div> <div>29% 71%</div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 855 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*AP*CP*CP*CP*GP*AP*CP*GP*GP*GP*AP*CP*TP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	21	Total	C	N	O	P	0	0	0
			430	203	88	119	20			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	P	0	0	0
			121	57	21	37	6			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	P	0	0	0
			160	78	27	48	7			

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

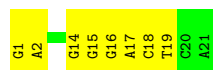
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	7	Total	C	N	O	P	0	0	0
			144	68	25	44	7			

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.

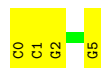
- Molecule 1: DNA (5'-D(*GP*AP*GP*CP*AP*GP*AP*CP*CP*CP*GP*AP*CP*GP*GP*GP*AP*CP*TP*CP*A)-3')

Chain A:  62% 38%



- Molecule 2: DNA (5'-D(P*CP*CP*GP*TP*CP*G)-3')

Chain B:  33% 67%




- Molecule 3: DNA (5'-D(*TP*CP*TP*GP*AP*GP*TP*C)-3')

Chain C:  62% 38%



- Molecule 4: DNA (5'-D(P*GP*GP*TP*CP*TP*GP*C)-3')

Chain D:  29% 71%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	68.17Å 68.17Å 53.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.08 – 3.16 34.08 – 3.16	Depositor EDS
% Data completeness (in resolution range)	66.3 (34.08-3.16) 73.5 (34.08-3.16)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.30 (at 3.18Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.201 , 0.216 0.206 , 0.225	Depositor DCC
R_{free} test set	322 reflections (9.26%)	wwPDB-VP
Wilson B-factor (Å ²)	152.5	Xtriage
Anisotropy	3.957	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.05 , -110.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.064 for -h,-k,l 0.076 for h,-h-k,-l 0.077 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.32	EDS
Total number of atoms	855	wwPDB-VP
Average B, all atoms (Å ²)	163.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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.

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	0.36	0/484	0.83	0/745
2	B	0.63	0/134	0.84	0/204
3	C	0.38	0/178	0.77	0/273
4	D	0.33	0/160	0.83	0/245
All	All	0.41	0/956	0.82	0/1467

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	430	0	234	7	0
2	B	121	0	68	5	1
3	C	160	0	93	2	1
4	D	144	0	80	6	1
All	All	855	0	475	17	2

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 13.

All (17) 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:5:DG:H2''	4:D:10:DG:O4'	2.00	0.61
3:C:1:DT:H2'	3:C:2:DC:C6	2.40	0.56
1:A:15:DG:H2''	1:A:16:DG:OP2	2.09	0.52
1:A:1:DG:H2''	1:A:2:DA:O5'	2.11	0.49
1:A:18:DC:H2''	1:A:19:DT:OP2	2.11	0.49
2:B:0:DC:H2'	2:B:1:DC:C6	2.48	0.49
2:B:5:DG:C2'	4:D:10:DG:O4'	2.60	0.49
1:A:14:DG:C6	1:A:15:DG:C6	3.01	0.49
4:D:14:DT:H2''	4:D:15:DG:C8	2.49	0.47
3:C:1:DT:H2''	3:C:2:DC:O5'	2.15	0.46
1:A:16:DG:H2''	1:A:17:DA:C8	2.52	0.45
2:B:5:DG:C6	4:D:10:DG:C5	3.06	0.44
4:D:11:DG:H2'	4:D:12:DT:C6	2.52	0.44
2:B:1:DC:H2''	2:B:2:DG:H8	1.85	0.41
4:D:10:DG:H2'	4:D:11:DG:O4'	2.21	0.40
1:A:1:DG:H2'	1:A:2:DA:C8	2.57	0.40
1:A:17:DA:C4	1:A:18:DC:C5	3.10	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:0:DC:OP2	2:B:5:DG:O3'[3_655]	2.12	0.08
3:C:8:DC:O3'	4:D:10:DG:OP1[3_655]	2.13	0.07

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

6.1 Protein, DNA and RNA chains

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

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

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

6.4 Ligands

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

6.5 Other polymers

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