Full wwPDB X-ray Structure Validation Report (i)

May 18, 2020 – 07:33 am BST

PDB ID	:	1DCV
Title	:	B-DNA DECAMER WITH CENTRAL TA DINUCLEOTIDE
Authors	:	Eichman, B.F.; Vargason, J.M.; Mooers, B.H.M.; Ho, P.S.
Deposited on	:	1999-11-05
Resolution	:	2.50 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

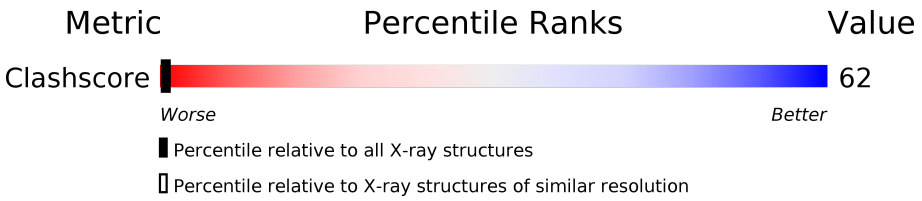
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.50 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))
Clashscore	141614	5346 (2.50-2.50)

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 >=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	10	10% 90%
1	B	10	100%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 427 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(*CP*CP*GP*CP*TP*AP*GP*CP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	A	10	202	96	39	58	9	0	0	0
1	B	10	202	96	39	58	9	0	0	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	13	13	13	0	0
2	B	10	10	10	0	0

3 Residue-property plots (i)

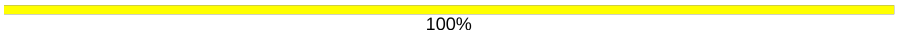
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according 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.

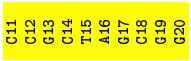
- Molecule 1: DNA (5'-D(*CP*CP*GP*CP*TP*AP*GP*CP*GP*G)-3')

Chain A:

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

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	64.10Å 25.90Å 39.90Å 90.00° 122.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.50 33.84 – 2.49	Depositor EDS
% Data completeness (in resolution range)	98.6 (10.00-2.50) 98.1 (33.84-2.49)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.48 (at 2.48Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.207 , 0.317 0.289 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	51.7	Xtriage
Anisotropy	1.047	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.23 , 878.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.42$, $< L^2 > = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	427	wwPDB-VP
Average B, all atoms (Å^2)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.33 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.8250e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] 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.

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	0.50	0/226	0.75	0/347
1	B	0.52	0/226	0.77	0/347
All	All	0.51	0/452	0.76	0/694

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	202	0	113	15	0
1	B	202	0	113	24	0
2	A	13	0	0	6	0
2	B	10	0	0	3	0
All	All	427	0	226	39	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 62.

All (39) 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:B:17:DG:H2"	1:B:18:DC:H5"	1.21	1.16

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:DG:C2'	1:B:18:DC:H5"	1.92	0.99
1:A:9:DG:H2"	1:A:10:DG:C8	2.10	0.86
1:B:19:DG:H2"	1:B:20:DG:C8	2.19	0.77
1:B:18:DC:H2'	1:B:19:DG:C8	2.23	0.74
1:A:1:DC:H2'	1:A:2:DC:C5	2.22	0.73
1:A:7:DG:N3	2:A:42:HOH:O	2.21	0.72
1:B:17:DG:H2"	1:B:18:DC:C5'	2.14	0.70
1:A:6:DA:H2	2:A:41:HOH:O	1.76	0.68
1:A:3:DG:H2"	1:A:4:DC:H5'	1.79	0.65
1:B:11:DC:O5'	1:B:11:DC:C6	2.48	0.64
1:B:17:DG:C3'	1:B:18:DC:H5"	2.28	0.64
1:A:6:DA:C2	2:A:41:HOH:O	2.51	0.61
1:A:8:DC:H2"	1:A:9:DG:C8	2.35	0.61
1:B:19:DG:H2"	1:B:20:DG:N7	2.17	0.59
1:B:17:DG:H2"	1:B:18:DC:C6	2.40	0.57
1:A:8:DC:O4'	2:A:42:HOH:O	2.18	0.56
1:A:7:DG:H2"	1:A:8:DC:OP2	2.07	0.55
1:A:1:DC:H5	2:A:30:HOH:O	1.89	0.55
1:B:18:DC:H6	1:B:18:DC:C5'	2.20	0.54
1:B:15:DT:H2"	1:B:16:DA:N7	2.25	0.51
1:A:1:DC:H6	1:A:1:DC:HO5'	1.58	0.50
1:B:19:DG:C2'	1:B:20:DG:N7	2.74	0.49
1:B:13:DG:H2"	1:B:14:DC:OP2	2.11	0.49
1:B:18:DC:C2'	1:B:19:DG:C8	2.94	0.48
1:A:3:DG:H1'	1:A:4:DC:H5"	1.95	0.48
1:B:15:DT:H2"	1:B:16:DA:C8	2.48	0.48
1:B:11:DC:H6	1:B:11:DC:HO5'	1.48	0.47
1:B:11:DC:H2'	1:B:12:DC:C5	2.50	0.46
1:A:1:DC:H2'	1:A:2:DC:C6	2.52	0.45
1:A:1:DC:H2'	1:A:2:DC:C4	2.52	0.45
1:B:18:DC:H5'	1:B:18:DC:H6	1.82	0.45
1:B:19:DG:H3'	2:B:39:HOH:O	2.18	0.44
1:B:19:DG:H5'	2:B:39:HOH:O	2.18	0.42
1:B:18:DC:H2'	1:B:19:DG:N7	2.33	0.42
1:A:7:DG:N2	2:A:25:HOH:O	2.44	0.41
1:B:11:DC:H6	1:B:11:DC:O5'	1.96	0.40
1:B:19:DG:C5'	2:B:39:HOH:O	2.69	0.40
1:B:20:DG:H2'	1:B:20:DG:O5'	2.21	0.40

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 carbohydrates 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)

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6.4 Ligands (i)

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

6.5 Other polymers (i)

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