



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 01:02 pm BST

PDB ID : 441D
Title : HIGH RESOLUTION A-DNA CRYSTAL STRUCTURES OF
D(AGGGGCCCT): AN A-DNA MODEL OF POLY(DG).POLY(DC)
Authors : Gao, Y.G.; Robinson, H.H.; Wang, A.H.
Deposited on : 1999-01-13
Resolution : 1.50 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 ⓘ symbol.

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	:	2.11
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

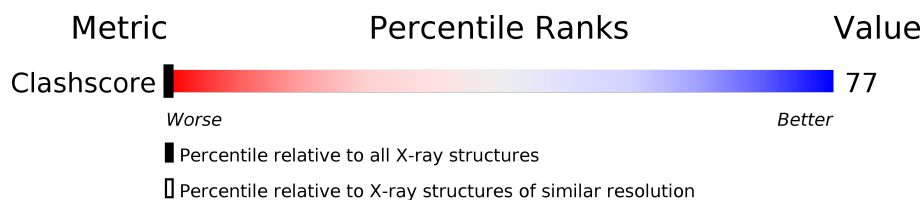
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.50 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	3144 (1.50-1.50)

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	10	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 280 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(*AP*GP*GP*GP*GP*CP*CP*CP*CP*T)-3').

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

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	78	Total	O	0	0
			78	78		

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(*AP*GP*GP*GP*GP*CP*CP*CP*CP*T)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	32.77Å 32.77Å 77.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 1.50 28.38 – 1.34	Depositor EDS
% Data completeness (in resolution range)	61.6 (6.00-1.50) 95.9 (28.38-1.34)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 1.34Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.236 , 0.315 0.276 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	7.1	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.50 , 262.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	280	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.49% 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

5.1 Standard geometry

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	2.11	10/226 (4.4%)	1.45	4/347 (1.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	3	DG	C2-N2	-8.96	1.25	1.34
1	A	6	DC	N1-C2	-7.68	1.32	1.40
1	A	3	DG	N1-C2	-7.23	1.31	1.37
1	A	1	DA	C6-N1	-7.01	1.30	1.35
1	A	6	DC	C4-C5	-6.41	1.37	1.43
1	A	4	DG	N9-C4	-6.24	1.32	1.38
1	A	4	DG	C3'-O3'	-5.64	1.36	1.44
1	A	8	DC	N1-C2	-5.24	1.34	1.40
1	A	3	DG	C6-N1	-5.17	1.35	1.39
1	A	1	DA	C6-N6	-5.02	1.29	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	5	DG	C5'-C4'-O4'	6.11	120.91	109.30
1	A	3	DG	C5'-C4'-O4'	5.67	120.07	109.30
1	A	1	DA	N1-C6-N6	-5.52	115.28	118.60
1	A	3	DG	N1-C2-N2	-5.37	111.37	116.20

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	3	DG	Sidechain
1	A	5	DG	Sidechain
1	A	6	DC	Sidechain
1	A	9	DC	Sidechain

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	202	0	110	24	0
2	A	78	0	0	4	0
All	All	280	0	110	24	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 77.

All (24) 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:7:DC:H6	1:A:7:DC:H5'	1.05	1.09
1:A:8:DC:H2''	1:A:9:DC:H5'	1.31	1.08
1:A:8:DC:H3'	2:A:104:HOH:O	1.56	1.05
1:A:7:DC:H5'	1:A:7:DC:C6	1.94	1.03
1:A:8:DC:C2'	1:A:9:DC:H5'	1.95	0.96
1:A:7:DC:C5'	1:A:7:DC:H6	1.86	0.88
1:A:8:DC:H2''	1:A:9:DC:C5'	2.04	0.86
1:A:8:DC:C2'	1:A:9:DC:C5'	2.62	0.74
1:A:6:DC:H2'	1:A:7:DC:H5''	1.72	0.72
1:A:7:DC:C5'	1:A:7:DC:C6	2.71	0.66
1:A:8:DC:H2'	1:A:9:DC:H5'	1.86	0.57
1:A:7:DC:H2'	1:A:8:DC:C6	2.39	0.57
1:A:6:DC:C2'	1:A:7:DC:H5''	2.35	0.55
1:A:10:DT:H2''	2:A:131:HOH:O	2.07	0.54
1:A:8:DC:H2'	1:A:9:DC:C5'	2.39	0.53
1:A:8:DC:C2'	1:A:9:DC:O5'	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:DG:P	2:A:118:HOH:O	2.73	0.46
1:A:7:DC:C4'	1:A:7:DC:C6	2.99	0.45
1:A:1:DA:H5''	1:A:2:DG:C8	2.53	0.44
1:A:8:DC:H6	1:A:8:DC:O5'	2.01	0.44
1:A:7:DC:C3'	1:A:7:DC:C6	3.02	0.43
1:A:8:DC:H2'	1:A:9:DC:O5'	2.19	0.43
1:A:6:DC:H5	2:A:173:HOH:O	2.01	0.43
1:A:3:DG:H2'	1:A:4:DG:O5'	2.20	0.42

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

There are no such residues in this entry.

5.8 Polymer linkage issues

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.