



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 03:00 PM GMT

PDB ID : 117D  
Title : CRYSTAL AND MOLECULAR STRUCTURE OF THE ALTERNATING  
DODECAMER D(GCGTACGTACGC) IN THE A-DNA FORM: COMPAR-  
ISON WITH THE ISOMORPHOUS NON-ALTERNATING DODECAMER  
D(CCGTACGTACGG)  
Authors : Bingman, C.; Jain, S.; Zon, G.; Sundaralingam, M.  
Deposited on : 1993-02-10  
Resolution : 2.55 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

---

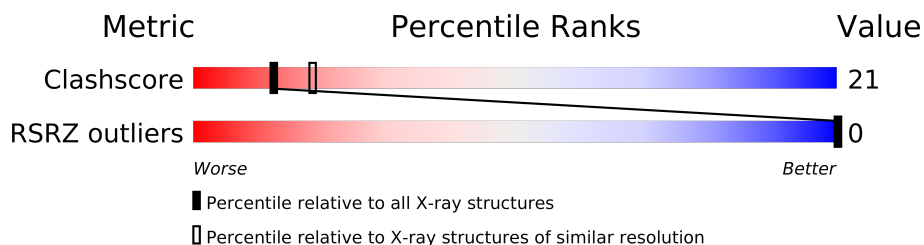
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.55 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	79885	4284 (2.58-2.50)
RSRZ outliers	66119	3414 (2.58-2.50)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	12	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 276 atoms, of which 0 are hydrogen and 0 are deuterium.

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\*CP\*GP\*TP\*AP\*CP\*GP\*TP\*AP\*CP\*GP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	12	Total	C	N	O	P	0	0	0
			243	116	46	70	11			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	33	Total	O	0	0
			33	33		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). 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\*CP\*GP\*TP\*AP\*CP\*GP\*TP\*AP\*CP\*GP\*C)-3')

Chain A: 

G1	C2	G3	T4	A5	C6	G7	T8	A9	C10	G11	C12
----	----	----	----	----	----	----	----	----	-----	-----	-----

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.20Å 46.20Å 71.50Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.55 26.66 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.55) 71.8 (26.66-2.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.57 (at 2.60Å)	Xtriage
Refinement program	NUCLSQ	Depositor
R, $R_{free}$	0.142 , (Not available) 0.153 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	38.7	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 67.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 1146 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	276	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.31% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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	1.54	1/272 (0.4%)	2.59	30/418 (7.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	3	DG	O3'-P	5.51	1.67	1.61

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	DG	O5'-P-OP1	10.11	122.83	110.70
1	A	2	DC	P-O3'-C3'	-10.09	107.59	119.70
1	A	8	DT	C2-N3-C4	-9.12	121.73	127.20
1	A	9	DA	C6-N1-C2	8.14	123.48	118.60
1	A	9	DA	N1-C2-N3	-7.88	125.36	129.30
1	A	8	DT	N3-C4-O4	-7.47	115.42	119.90
1	A	6	DC	O4'-C1'-N1	7.31	113.12	108.00
1	A	8	DT	N3-C4-C5	7.25	119.55	115.20
1	A	12	DC	O4'-C1'-N1	7.19	113.03	108.00
1	A	2	DC	P-O5'-C5'	-6.95	109.79	120.90
1	A	1	DG	P-O3'-C3'	-6.92	111.39	119.70
1	A	9	DA	C5-C6-N1	-6.83	114.28	117.70
1	A	8	DT	O5'-P-OP1	6.72	118.77	110.70
1	A	5	DA	O4'-C1'-N9	6.51	112.56	108.00
1	A	11	DG	C5-C6-N1	6.28	114.64	111.50
1	A	3	DG	C5-C6-N1	6.23	114.62	111.50
1	A	4	DT	C2-N3-C4	-6.12	123.53	127.20
1	A	6	DC	N3-C4-C5	-5.85	119.56	121.90
1	A	1	DG	C5-C6-N1	5.74	114.37	111.50
1	A	7	DG	OP1-P-OP2	-5.59	111.22	119.60
1	A	1	DG	O4'-C1'-N9	5.52	111.87	108.00
1	A	6	DC	N1-C2-O2	5.47	122.18	118.90
1	A	8	DT	N1-C2-N3	5.45	117.87	114.60

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	DA	OP1-P-OP2	-5.38	111.53	119.60
1	A	5	DA	N1-C2-N3	-5.26	126.67	129.30
1	A	5	DA	O5'-P-OP1	5.25	117.00	110.70
1	A	7	DG	O5'-P-OP1	5.21	116.95	110.70
1	A	12	DC	N3-C4-C5	-5.09	119.86	121.90
1	A	12	DC	OP1-P-OP2	-5.08	111.98	119.60
1	A	11	DG	OP2-P-O3'	5.02	116.23	105.20

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	243	0	136	8	1
2	A	33	0	0	5	1
All	All	276	0	136	8	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 21.

All (8) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:9:DA:H2''	2:A:42:HOH:O	1.89	0.71
1:A:8:DT:H1'	2:A:23:HOH:O	2.02	0.59
1:A:1:DG:H5''	2:A:14:HOH:O	2.07	0.53
1:A:9:DA:C2'	2:A:42:HOH:O	2.55	0.51
1:A:8:DT:H2''	1:A:9:DA:H5'	1.95	0.47
1:A:9:DA:C2'	1:A:10:DC:O5'	2.63	0.46
1:A:1:DG:H5''	1:A:2:DC:OP2	2.16	0.46
1:A:9:DA:H1'	2:A:42:HOH:O	2.16	0.45

All (1) 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	Distance(Å)	Clash(Å)
1:A:12:DC:O3'	2:A:23:HOH:O[5_554]	2.17	0.03

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

There are no protein chains in this entry.

### 5.3.2 Protein sidechains ⓘ

There are no protein chains in this entry.

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	12/12 (100%)	-0.84	0 <span>100</span> <span>100</span>	16, 21, 25, 32	0

There are no RSRZ outliers to report.

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

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates

There are no carbohydrates in this entry.

### 6.4 Ligands

There are no ligands in this entry.

### 6.5 Other polymers

There are no such residues in this entry.