



wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 1D68 / pdb_00001d68
Title : SOLUTION STRUCTURE OF [D(GCGTATACGC)]2
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Deposited on : 1992-04-15

This is a wwPDB NMR 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/NMRValidationReportHelp>

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 : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Ensemble composition and analysis

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 636 atoms, of which 224 are hydrogens and 0 are deuteriums.


- Molecule 1 is a DNA chain called DNA (5'-D(P*GP*CP*GP*TP*AP*TP*AP*CP*GP*C)-3').

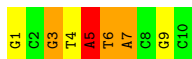
| Mol | Chain | Residues | Atoms | | | | | Trace | |
|-----|-------|----------|-------|----|-----|----|----|-------|---|
| | | | Total | C | H | N | O | | P |
| 1 | A | 10 | 318 | 97 | 112 | 38 | 61 | 10 | 0 |
| 1 | B | 10 | 318 | 97 | 112 | 38 | 61 | 10 | 0 |

4 Residue-property plots [i](#)


These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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 outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

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

Chain A: 



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

Chain B: 



5 Refinement protocol and experimental data overview

Of the ? calculated structures, 1 were deposited, based on the following criterion: ?.

The authors did not provide any information on software used for structure solution, optimization or refinement.

No chemical shift data was provided.

6 Model quality i

6.1 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 (average) 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 | 0.97 | 1/230 (0.4%) | 1.48 | 5/351 (1.4%) |
| 1 | B | 0.97 | 1/230 (0.4%) | 1.47 | 5/351 (1.4%) |
| All | All | 0.97 | 2/460 (0.4%) | 1.47 | 10/702 (1.4%) |

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 | Planarity |
|-----|-------|-----------|-----------|
| 1 | A | 0 | 2 |
| 1 | B | 0 | 2 |
| All | All | 0 | 4 |

All bond outliers are listed below. They are sorted according to the Z-score.

| Mol | Chain | Res | Type | Atoms | Z | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|-------|------|-------------|----------|
| 1 | B | 11 | DG | P-OP2 | 5.99 | 1.60 | 1.48 |
| 1 | A | 1 | DG | P-OP2 | 5.99 | 1.60 | 1.48 |

5 of 10 angle outliers are listed below. They are sorted according to the Z-score.

| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|-----|------|-----------|------|-------------|----------|
| 1 | A | 7 | DA | P-O3'-C3' | 8.21 | 132.52 | 120.20 |
| 1 | B | 17 | DA | P-O3'-C3' | 8.20 | 132.50 | 120.20 |
| 1 | A | 5 | DA | P-O3'-C3' | 8.00 | 132.21 | 120.20 |
| 1 | B | 15 | DA | P-O3'-C3' | 7.98 | 132.17 | 120.20 |
| 1 | B | 11 | DG | P-O3'-C3' | 7.70 | 131.75 | 120.20 |

There are no chirality outliers.

All planar outliers are listed below.

| Mol | Chain | Res | Type | Group |
|-----|-------|-----|------|-----------|
| 1 | A | 5 | DA | Sidechain |
| 1 | A | 9 | DG | Sidechain |
| 1 | B | 15 | DA | Sidechain |
| 1 | B | 19 | DG | Sidechain |

6.2 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 each 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 averaged over the ensemble.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes |
|-----|-------|-------|----------|----------|---------|
| 1 | A | 206 | 112 | 113 | 11 |
| 1 | B | 206 | 112 | 113 | 11 |
| All | All | 412 | 224 | 226 | 19 |

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 30.

5 of 19 clashes are listed below, sorted by their clash magnitude.

| Atom-1 | Atom-2 | Clash(Å) | Distance(Å) |
|--------------|--------------|----------|-------------|
| 1:A:5:DA:N6 | 1:B:15:DA:N6 | 0.55 | 2.55 |
| 1:A:7:DA:C2 | 1:B:15:DA:C2 | 0.54 | 2.95 |
| 1:B:14:DT:C4 | 1:B:15:DA:N6 | 0.54 | 2.76 |
| 1:A:4:DT:C4 | 1:A:5:DA:N6 | 0.54 | 2.76 |
| 1:A:5:DA:C2 | 1:B:17:DA:C2 | 0.53 | 2.96 |

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

There are no protein molecules in this entry.

6.3.2 Protein sidechains [i](#)

There are no protein molecules in this entry.

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided