



Full wwPDB NMR Structure Validation Report ⓘ

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PDB ID : 1N2W
Title : Solution Structure of 8OG:G mismatch containing duplex
Authors : Thiviyathan, V.; Somasunderam, A.D.; Hazra, T.K.; Mitra, S.; Gorenstein, D.G.
Deposited on : 2002-10-24

This is a Full wwPDB NMR 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/NMRValidationReportHelp>

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:

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : 2.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

2 Ensemble composition and analysis

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

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 766 atoms, of which 272 are hydrogens and 0 are deuteriums.

- Molecule 1 is a DNA chain called 5'-D>(*CP*GP*CP*GP*AP*AP*TP*TP*(8OG)P*GP*CP*G)-3'.

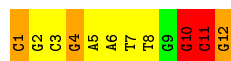
Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		P
1	A	12	Total 383	C 117	H 136	N 48	O 71	P 11	0
1	B	12	Total 383	C 117	H 136	N 48	O 71	P 11	0

4 Residue-property plots [i](#)

These plots are provided for all protein, RNA and DNA 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.

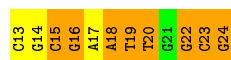
- Molecule 1: 5'-D(*CP*GP*CP*GP*AP*AP*TP*TP*(8OG)P*GP*CP*G)-3'

Chain A: 



- Molecule 1: 5'-D(*CP*GP*CP*GP*AP*AP*TP*TP*(8OG)P*GP*CP*G)-3'

Chain B: 



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *Relaxation matrix refinement, molecular dynamics, simulated annealing.*

Of the 1 calculated structures, 1 were deposited, based on the following criterion: *all calculated structures submitted.*

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
Amber	refinement	5.0

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 6 of this report.

Chemical shift file(s)	input_cs.cif
Number of chemical shift lists	1
Total number of shifts	232
Number of shifts mapped to atoms	218
Number of unparsed shifts	0
Number of shifts with mapping errors	14
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	47%

No validations of the models with respect to experimental NMR restraints is performed at this time.

COVALENT-GEOMETRY INFOmissingINFO

5.1 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	247	136	132	1
All	All	494	272	261	1

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 1.

All clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:10:DG:N2	1:A:11:DC:C2	0.41	2.89

5.2 Torsion angles [i](#)

5.2.1 Protein backbone [i](#)

There are no protein molecules in this entry.

5.2.2 Protein sidechains [i](#)

There are no protein molecules in this entry.

5.2.3 RNA [i](#)

There are no RNA molecules in this entry.

MODRES-GEOMETRY INFOmissingINFO

5.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.4 Ligand geometry [i](#)

There are no ligands in this entry.

5.5 Other polymers [i](#)

There are no such molecules in this entry.

5.6 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 47% for the well-defined parts and 47% for the entire structure.

6.1 Chemical shift list 1

File name: input_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

6.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	232
Number of shifts mapped to atoms	218
Number of unparsed shifts	0
Number of shifts with mapping errors	14
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- Residue not found in structure. All 14 occurrences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
A	9	DG	H1'	5.17	0.1	1
B	21	DG	H4'	4.11	0.1	1
B	21	DG	H2''	2.34	0.1	1
A	9	DG	H2'	3.27	0.1	1
A	9	DG	H4'	4.11	0.1	1
A	9	DG	H3'	4.95	0.1	1
B	21	DG	H1'	5.17	0.1	1
B	21	DG	H2'	3.27	0.1	1
B	21	DG	H5'	4.09	0.1	1
A	9	DG	H5'	4.09	0.1	1
B	21	DG	H3'	4.95	0.1	1
B	21	DG	H1	11.38	0.1	1
A	9	DG	H2''	2.34	0.1	1
A	9	DG	H1	11.38	0.1	1

6.1.2 Chemical shift referencing [i](#)

No chemical shift referencing corrections were calculated (not enough data).

6.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 47%, i.e. 208 atoms were assigned a chemical shift out of a possible 440. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	208/440 (47%)	208/264 (79%)	0/146 (0%)	0/30 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 47%, i.e. 208 atoms were assigned a chemical shift out of a possible 440. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	208/440 (47%)	208/264 (79%)	0/146 (0%)	0/30 (0%)

6.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

6.1.5 Random Coil Index (RCI) plots [i](#)

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned_chem_shift_list_1). RCI is only applicable to proteins.