



Full wwPDB NMR Structure Validation Report ⓘ

Mar 19, 2026 – 08:25 PM UTC

PDB ID : 1D6D / pdb_00001d6d
Title : SOLUTION DNA STRUCTURE CONTAINING (A-A)-T TRIADS INTER-DIGITATED BETWEEN A-T BASE PAIRS AND GGGG TETRADS; NMR, 8 STRUCT.
Authors : Kuryavyi, V.V.; Kettani, A.; Wang, W.; Jones, R.; Patel, D.J.
Deposited on : 1999-10-13

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 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

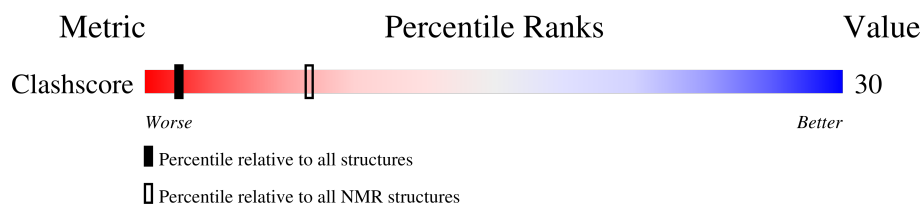
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

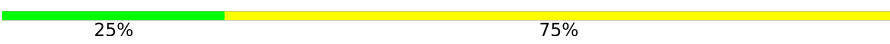
The overall completeness of chemical shifts assignment was not calculated.

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)	NMR archive (#Entries)
Clashscore	229148	14424

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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	12	 25% 75%
1	B	12	 25% 75%

2 Ensemble composition and analysis ⓘ

This entry contains 8 models. This entry does not contain polypeptide chains, therefore identification of well-defined residues and clustering analysis are not possible. All residues are included in the validation scores.

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 774 atoms, of which 276 are hydrogens and 0 are deuteriums.

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

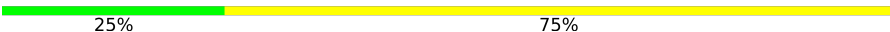
Mol	Chain	Residues	Atoms						Trace
1	A	12	Total	C	H	N	O	P	0
			387	120	138	48	70	11	
1	B	12	Total	C	H	N	O	P	0
			387	120	138	48	70	11	

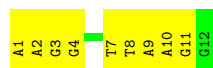
4 Residue-property plots [i](#)

4.1 Average score per residue in the NMR ensemble

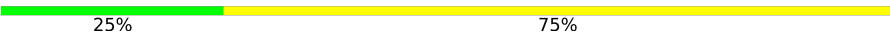
These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for 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.

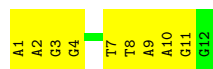
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

Chain A: 



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

Chain B: 

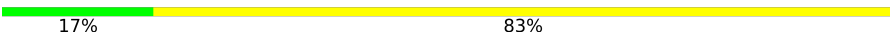


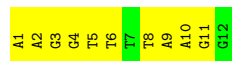
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

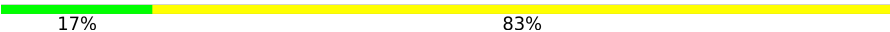
4.2.1 Score per residue for model 1

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

Chain A: 



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

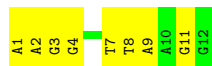
Chain B: 



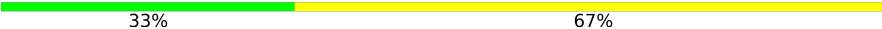
4.2.2 Score per residue for model 2

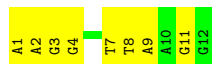
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

Chain A:  33% 67%



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

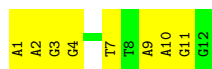
Chain B:  33% 67%



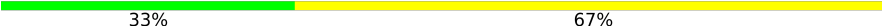
4.2.3 Score per residue for model 3

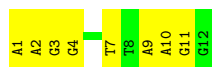
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

Chain A:  33% 67%



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

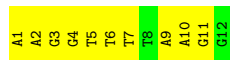
Chain B:  33% 67%



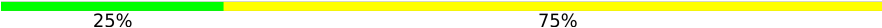
4.2.4 Score per residue for model 4

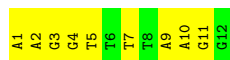
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

Chain A:  17% 83%



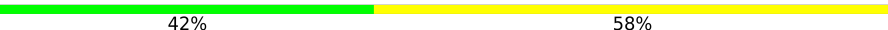
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

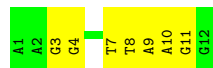
Chain B:  25% 75%



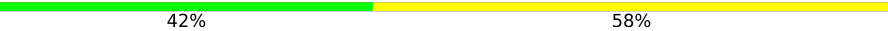
4.2.5 Score per residue for model 5

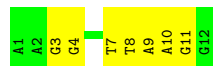
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

Chain A:  42% 58%



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

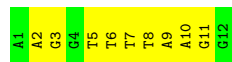
Chain B:  42% 58%



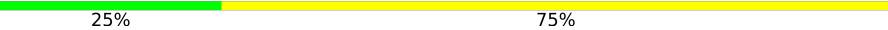
4.2.6 Score per residue for model 6

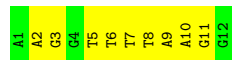
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

Chain A:  25% 75%



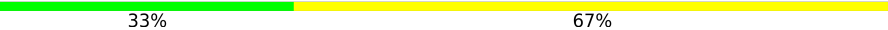
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

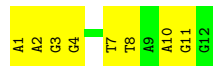
Chain B:  25% 75%



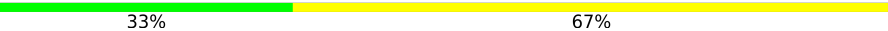
4.2.7 Score per residue for model 7

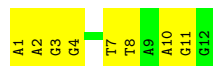
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

Chain A:  33% 67%



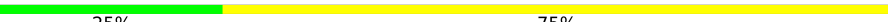
- Molecule 1: 5'-D(*AP*AP*GP*GP*TP*TP*TP*TP*AP*AP*GP*G)-3'

Chain B:  33% 67%



4.2.8 Score per residue for model 8

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

Chain A:  25% 75%



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

Chain B:  33% 67%



5 Refinement protocol and experimental data overview

The models were refined using the following method: *DISTANCE GEOMETRY AND SIMULATED ANNEALING, DISTANCE RESTRAINED MOLECULAR DYNAMICS REFINEMENT AND RELAXATION MATRIX INTENSITY REFINEMENT*.

Of the 100 calculated structures, 8 were deposited, based on the following criterion: *BACK CALCULATED DATA AGREE WITH EXPERIMENTAL NOESY SPECTRUM, STRUCTURES WITH ACCEPTABLE COVALENT GEOMETRY, STRUCTURES WITH FAVORABLE NON-BOND ENERGY, STRUCTURES WITH THE LEAST RESTRAINT VIOLATIONS, STRUCTURES WITH THE LOWEST ENERGY*.

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

Software name	Classification	Version
X-PLOR	refinement	3.1
X-PLOR	structure solution	3.1

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	249	138	126	13±2
1	B	249	138	126	12±2
All	All	3984	2208	2016	179

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.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:9:DA:H1'	1:B:10:DA:N7	0.88	1.83	3	3
1:A:9:DA:H1'	1:A:10:DA:N7	0.87	1.83	3	3
1:A:6:DT:H3'	1:A:6:DT:OP2	0.73	1.84	1	1
1:B:6:DT:H3'	1:B:6:DT:OP2	0.72	1.84	1	1
1:B:8:DT:O3'	1:B:9:DA:H3'	0.71	1.85	2	2
1:A:8:DT:O3'	1:A:9:DA:H3'	0.71	1.85	2	2
1:A:1:DA:H2'	1:B:7:DT:O4	0.68	1.88	7	1
1:A:7:DT:O4	1:B:1:DA:H2'	0.68	1.88	7	1
1:B:10:DA:H4'	1:B:11:DG:C5'	0.65	2.22	7	1
1:A:10:DA:H4'	1:A:11:DG:C5'	0.65	2.22	7	1
1:A:1:DA:H3'	1:B:7:DT:O4	0.64	1.93	2	1
1:A:11:DG:H2'	1:A:11:DG:N3	0.63	2.08	1	7
1:A:7:DT:O4	1:B:1:DA:H3'	0.63	1.93	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:11:DG:H2'	1:B:11:DG:N3	0.62	2.08	1	4
1:A:3:DG:H2''	1:A:4:DG:O5'	0.61	1.96	1	7
1:B:3:DG:H2''	1:B:4:DG:O5'	0.60	1.96	1	7
1:A:10:DA:H4'	1:A:11:DG:O5'	0.60	1.97	3	5
1:B:10:DA:H4'	1:B:11:DG:O5'	0.60	1.97	3	5
1:A:1:DA:H2''	1:A:2:DA:O5'	0.59	1.97	3	4
1:B:1:DA:H2''	1:B:2:DA:O5'	0.59	1.97	3	4
1:A:7:DT:H4'	1:A:8:DT:OP1	0.58	1.98	7	2
1:B:7:DT:H4'	1:B:8:DT:OP1	0.57	1.98	7	2
1:B:5:DT:H2''	1:B:6:DT:OP2	0.56	2.01	1	3
1:B:11:DG:N3	1:B:11:DG:H2'	0.56	2.15	2	3
1:A:10:DA:N6	1:B:2:DA:C8	0.55	2.75	7	2
1:A:5:DT:H2''	1:A:6:DT:OP2	0.54	2.01	1	3
1:A:2:DA:C8	1:B:10:DA:N6	0.54	2.75	7	2
1:B:3:DG:N3	1:B:3:DG:H2'	0.54	2.16	8	4
1:A:3:DG:N3	1:A:3:DG:H2'	0.54	2.16	8	8
1:B:7:DT:H1'	1:B:9:DA:C5	0.51	2.41	5	1
1:A:7:DT:H1'	1:A:9:DA:C5	0.51	2.41	5	1
1:B:3:DG:H2'	1:B:3:DG:N3	0.51	2.20	1	4
1:A:9:DA:H2''	1:B:2:DA:N6	0.50	2.22	4	4
1:A:2:DA:N6	1:B:9:DA:H2''	0.50	2.22	4	4
1:A:10:DA:H1'	1:A:11:DG:C2	0.50	2.42	3	3
1:B:1:DA:N7	1:B:2:DA:C6	0.50	2.80	4	1
1:B:10:DA:H1'	1:B:11:DG:C2	0.49	2.42	3	3
1:A:1:DA:N7	1:A:2:DA:C6	0.49	2.80	4	1
1:B:7:DT:O2	1:B:9:DA:C8	0.49	2.65	5	1
1:A:10:DA:H1'	1:A:11:DG:N2	0.48	2.23	5	2
1:B:10:DA:H1'	1:B:11:DG:N2	0.48	2.23	5	2
1:B:7:DT:O4'	1:B:9:DA:C6	0.48	2.67	3	2
1:A:7:DT:O2	1:A:9:DA:C8	0.48	2.65	5	1
1:A:2:DA:H61	1:B:9:DA:H2''	0.48	1.69	2	1
1:B:10:DA:H4'	1:B:11:DG:H5'	0.47	1.86	7	1
1:B:5:DT:C2'	1:B:6:DT:OP2	0.47	2.63	6	2
1:A:10:DA:H4'	1:A:11:DG:H5'	0.47	1.86	7	1
1:A:7:DT:O4'	1:A:9:DA:C6	0.47	2.67	3	2
1:A:9:DA:H2''	1:B:2:DA:H61	0.46	1.69	2	1
1:B:10:DA:H4'	1:B:11:DG:H5''	0.46	1.86	7	1
1:A:7:DT:OP2	1:A:7:DT:H71	0.46	2.11	5	1
1:A:5:DT:C2'	1:A:6:DT:OP2	0.45	2.63	6	2
1:B:7:DT:H71	1:B:7:DT:OP2	0.45	2.11	5	1
1:A:8:DT:OP1	1:A:9:DA:C2	0.45	2.70	5	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:8:DT:OP1	1:B:9:DA:C2	0.44	2.70	5	1
1:A:3:DG:H1'	1:A:4:DG:O4'	0.44	2.13	4	1
1:B:1:DA:C8	1:B:2:DA:C4	0.43	3.06	1	2
1:A:10:DA:H4'	1:A:11:DG:H5''	0.43	1.87	7	1
1:B:3:DG:H1'	1:B:4:DG:O4'	0.43	2.13	4	1
1:A:1:DA:C8	1:A:2:DA:C4	0.43	3.06	1	3
1:A:1:DA:N7	1:A:2:DA:C5	0.42	2.88	1	1
1:A:4:DG:H2''	1:A:5:DT:O5'	0.42	2.15	4	1
1:B:1:DA:N7	1:B:2:DA:C5	0.41	2.88	1	1
1:A:11:DG:N3	1:A:11:DG:H2'	0.41	2.30	6	1
1:A:10:DA:C6	1:B:2:DA:C8	0.41	3.08	7	1
1:B:4:DG:H2''	1:B:5:DT:O5'	0.41	2.15	4	1
1:A:10:DA:N1	1:B:2:DA:C4	0.40	2.90	6	1
1:A:2:DA:C4	1:B:10:DA:N1	0.40	2.90	6	1

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

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided