

Full wwPDB NMR Structure Validation Report (i)

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PDB ID	:	7D0Z
Title	:	NMR solution structures of the DNA minidumbbell formed by two CCTG
		repeats at pH 5
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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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

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	:	2.16
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)		
Validation Pipeline (wwPDB-VP)	:	2.16

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $SOLUTION\ NMR$

The overall completeness of chemical shifts assignment is 44%.

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	Percer	ntile Ranks	Value
Clashscore			0
	Worse		Better
	Percentile relative to all structu		
	Percentile relative to all NMR st	tructures	
Motrio	Whole archive	NMR archive	

Metric	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f NMR} { m archive} \ (\#{ m Entries})$
Clashscore	158937	12864

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 <=5%

Mol	Chain	Length		Quality of chain	
1	А	8	13%	63%	25%



2 Ensemble composition and analysis (i)

This entry contains 20 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 are 2 unique types of molecules in this entry. The entry contains 256 atoms, of which 94 are hydrogens and 0 are deuteriums.

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

Mol	Chain	Residues		ŀ	4ton	ns			Trace
1	٨	0	Total	С	Η	Ν	0	Р	0
	A	0	251	76	94	26	48	7	0

• Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	
2	А	5	Total Na 5 5	



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: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	13%	63%	25%
DCZ1 C2 C3 C5 C5 C6 C6 C6 C6 C6 C6 C6 C6 C6 C6 C6 C7 C7 C7 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2			

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: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A: 13% 38% 50%

4.2.2 Score per residue for model 2

Chain A:	13%	50%	38%
DCZ1 C2 C3 C3 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5			



4.2.3 Score per residue for model 3

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	13%	63%	25%
DCZ1 C2 C3 64 C5 C5 C5 C5 C6 C5 C6 C6 C6 C5 C6 C6 C6 C7 C7 C7 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2			

4.2.4 Score per residue for model 4

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	75%	25%
DCZ1 C2 T3 G4 C5 C6 C6 C6 C6 C6 C6 C6 C6 C6 C6 C6 C6 C6		

4.2.5 Score per residue for model 5

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

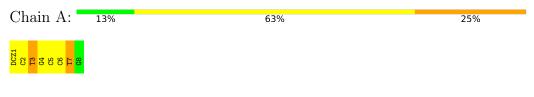
Chain A:	13%	63%	25%
DCZ1 C2 C3 64 64 C5 C5 C6 C5 C6 C6 C6 C6 C6 C6 C6 C7 C7 C7 C7 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2			

4.2.6 Score per residue for model 6

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	13%	63%	25%
DCZ1 C2 C3 G4 C5 C6 C6 C6 C6 C6 C6 C6 C6 C6 C6 C6 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7			

4.2.7 Score per residue for model 7





4.2.8 Score per residue for model 8

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	25%	50%	25%
DCZ1 C2 T3 G5 C5 C5 C5 C6 C5 C6 C5 C6 C6 C6 C5 C6 C6 C6 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7			

4.2.9 Score per residue for model 9

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	25%	38%	38%
DCZ1 C2 C3 C5 C5 C5 C5 C5 C5 C5 C6 C6 C6 C6 C6 C6 C6 C6 C7 C7 C7 C2 C7 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2			

4.2.10 Score per residue for model 10

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	25%	50%	25%
DCZ1 C2 13 64 64 65 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5			

4.2.11 Score per residue for model 11

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	63%	38%
DC2 C2 C3 C3 C5 C5 C5 C6 C6 C6 C6 C6 C6 C6 C6 C6 C7 C1 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2		

4.2.12 Score per residue for model 12

Chain A:	63%	38%
0 C Z 1 C 2 C 2 C 4 C 5 C 5 C 5 C 5 C 6 C 6 C 6 C 7 C 7 C 7 C 7 C 7 C 7 C 7 C 7 C 7 C 7		



4.2.13 Score per residue for model 13

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	13%	50%	38%
DC21 C2 C3 C4 C5 C5 C5 C6 C5 C6 C5 C6 C5 C6 C6 C6 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7			

4.2.14 Score per residue for model 14

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	13%	75%	13%
DCZ1 C2 C2 C5 C5 C5 C5 C6 C6 C6 C6 C6			

4.2.15 Score per residue for model 15

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

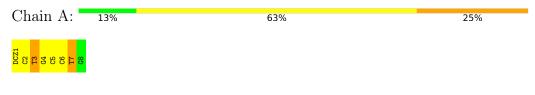
Chain A:	13%	50%	38%
DCZ1 C2 C3 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5 C5			

4.2.16 Score per residue for model 16

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	13%	63%	25%
DCZ1 C2 C3 C5 C5 C5 C6 C6 C5 C6 C6 C6 C6 C6 C6 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7 C7			

4.2.17 Score per residue for model 17





4.2.18 Score per residue for model 18

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	13%	75%	13%
DCZ1 C2 C2 C2 C5 C5 C5 C5 C5 C6 C6 C6 C6			

4.2.19 Score per residue for model 19

• Molecule 1: DNA (5'-D(*(DCZ)P*CP*TP*GP*CP*CP*TP*G)-3')

Chain A:	13%	63%	25%
DCZ1 C2 C3 C5 C5 C5 C5 C6 C5 C6 C5 C6 C5 C6 C5 C6 C5 C6 C5 C5 C5 C5 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2			

4.2.20 Score per residue for model 20

Chain A:	75%	25%
DCZ1 C2 C3 C4 C5 C5 C5 C6 C5 C6 C5 C6 C5 C6 C5 C6 C5 C6 C7 C7 C7 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2 C2		



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: restrained molecular dynamics.

Of the 1000 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
Amber	structure calculation	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	76
Number of shifts mapped to atoms	76
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	44%



6 Model quality (i)

6.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, DCZ

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	B	ond lengths	Bond angles		
	Chain	RMSZ $\#Z > 5$		RMSZ $\#Z>5$		
1	А	$1.38 {\pm} 0.03$	$0{\pm}0/156~(~0.0{\pm}~0.0\%)$	$2.32{\pm}0.06$	$12{\pm}2/238~(~5.1{\pm}~0.7\%)$	
All	All	1.38	0/3120~(~0.0%)	2.32	243/4760 ($5.1%$)	

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	А	$0.0{\pm}0.0$	2.5 ± 0.7
All	All	0	50

There are no bond-length outliers.

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mal	Mol Chain		Turne	Atoms	Z	Observed(°)	$Ideal(^{o})$	Mo	dels
	Unain	Res	Type	Atoms		Observed(*)	Ideal(*)	Worst	Total
1	А	7	DT	C6-C5-C7	-12.07	115.66	122.90	18	20
1	А	7	DT	O4'-C1'-N1	9.15	114.41	108.00	15	16
1	А	5	DC	O4'-C1'-N1	8.63	114.04	108.00	4	9
1	А	5	DC	C1'-O4'-C4'	-8.17	101.93	110.10	1	8
1	А	2	DC	O4'-C1'-N1	8.14	113.70	108.00	13	20
1	А	7	DT	P-O3'-C3'	8.05	129.36	119.70	20	1
1	А	4	DG	O4'-C1'-N9	7.94	113.56	108.00	8	11
1	А	6	DC	P-O3'-C3'	7.43	128.61	119.70	13	10
1	А	6	DC	O4'-C1'-N1	7.17	113.02	108.00	13	17
1	А	4	DG	N3-C4-C5	-6.97	125.12	128.60	6	19
1	А	7	DT	C4-C5-C6	6.48	121.89	118.00	5	20
1	А	3	DT	O4'-C1'-N1	6.44	112.51	108.00	13	3

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Mal	Chair	Dec	True	Atoma	7	Observed ⁽⁰⁾		Moo	dels
Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$\mathrm{Ideal}(^{o})$	Worst	Total
1	А	6	DC	N1-C2-O2	6.42	122.75	118.90	2	20
1	А	3	DT	P-O3'-C3'	6.41	127.39	119.70	4	9
1	А	8	DG	O4'-C1'-N9	6.21	112.35	108.00	16	4
1	А	5	DC	P-O3'-C3'	6.16	127.09	119.70	2	3
1	А	2	DC	P-O3'-C3'	6.11	127.03	119.70	17	1
1	А	7	DT	C5-C6-N1	-5.91	120.16	123.70	5	15
1	А	3	DT	C6-C5-C7	-5.89	119.37	122.90	15	18
1	А	5	DC	N1-C1'-C2'	5.43	122.91	112.60	1	4
1	А	6	DC	N3-C2-O2	-5.34	118.16	121.90	19	7
1	А	4	DG	C8-N9-C4	-5.29	104.28	106.40	6	2
1	А	7	DT	C4-C5-C7	5.25	122.15	119.00	11	2
1	А	8	DG	C8-N9-C4	-5.21	104.31	106.40	15	2
1	А	7	DT	N3-C2-O2	-5.08	119.25	122.30	5	2

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There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	А	7	DT	Sidechain	20
1	А	3	DT	Sidechain	13
1	А	4	DG	Sidechain	7
1	А	5	DC	Sidechain	4
1	А	8	DG	Sidechain	3
1	А	6	DC	Sidechain	2
1	А	2	DC	Sidechain	1

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
	All	All	3240	1880	1840	-

The all-atom clash score is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clash score for this structure is -.

There are no clashes.



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)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with |Z| > 2 is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mal	Tuno	Chain	Dog	Link		Bond leng	ths
	Type	Ullalli	nes		Counts	RMSZ	#Z>2
1	DCZ	А	1	1	$15,\!17,\!17$	$0.80{\pm}0.00$	$0\pm0~(0\pm0\%)$

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with |Z| > 2 is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mal	Turne	Chain	Dec	Tink		Bond ang	gles
WIOI	Type	Unam	nes	LIIIK	Counts	RMSZ	#Z>2
1	DCZ	А	1	1	17,24,24	$1.44{\pm}0.00$	$1\pm0 (5\pm0\%)$

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical



component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	DCZ	А	1	1	-	$0\pm 0,3,18,18$	$0\pm 0,2,2,2$

There are no bond-length outliers.

All unique angle outliers are listed below.

Mo	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$\operatorname{Ideal}(^{o})$	Moo Worst	
1	A	1	DCZ	C2-N3-C4	5.15	121.56	116.34	9	20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates (i)

There are no monosaccharides in this entry.

6.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

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 (i)

The completeness of assignment taking into account all chemical shift lists is 44% for the well-defined parts and 44% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: cctg2ph5_cs_deposit

7.1.1 Bookkeeping (i)

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	76
Number of shifts mapped to atoms	76
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing (i)

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

7.1.3 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 44%, i.e. 61 atoms were assigned a chemical shift out of a possible 139. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^{1}\mathbf{H}$	$^{13}\mathrm{C}$	$^{15}\mathbf{N}$
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	61/139~(44%)	61/82~(74%)	0/47~(0%)	0/10~(0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 44%, i.e. 61 atoms were assigned a chemical shift out of a possible 139. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.



	Total	$^{1}\mathbf{H}$	$^{13}\mathrm{C}$	$^{15}\mathbf{N}$
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	61/139~(44%)	61/82~(74%)	0/47~(0%)	0/10~(0%)

7.1.4 Statistically unusual chemical shifts (i)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots (i)

No random coil index (RCI) plot could be generated from the current chemical shift list (cctg2ph5_cs_deposit) RCI is only applicable to proteins.

