



wwPDB NMR Structure Validation Summary Report

Nov 20, 2023 – 10:19 AM EST


PDB ID : 2O7X
Title : Duplex DNA containing an abasic site with an opposite G (beta anomer) in 5'-G_AC-3' (10 structure ensemble and averaged structure)
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Deposited on : 2006-12-11

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

<https://www.wwpdb.org/validation/2017/FAQs#types>.

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)
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) : 2.36

2 Ensemble composition and analysis

This entry contains 11 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 811 atoms, of which 291 are hydrogens and 0 are deuteriums.

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

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
1	A	13	399	121	142	52	72	12	0

- Molecule 2 is a DNA chain called 5'-D(*CP*CP*CP*GP*GP*TP*GP*CP*TP*TP*TP*GP*G)-3'.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
2	B	13	412	126	149	45	80	12	0

4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

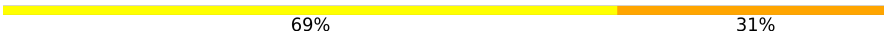
These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for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

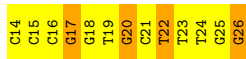
- Molecule 1: 5'-D(*CP*CP*AP*AP*AP*GP*(AAB)P*AP*CP*CP*GP*GP*G)-3'

Chain A: 



- Molecule 2: 5'-D(*CP*CP*CP*GP*GP*TP*GP*CP*TP*TP*TP*GP*G)-3'

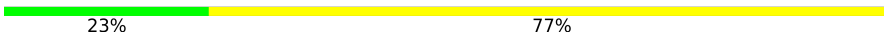
Chain B: 

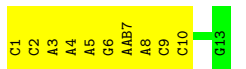


4.2 Residue scores for the representative (author defined) model from the NMR ensemble

The representative model is number 11. Colouring as in section 4.1 above.

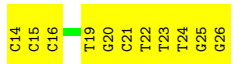
- Molecule 1: 5'-D(*CP*CP*AP*AP*AP*GP*(AAB)P*AP*CP*CP*GP*GP*G)-3'

Chain A: 



- Molecule 2: 5'-D(*CP*CP*CP*GP*GP*TP*GP*CP*TP*TP*TP*GP*G)-3'

Chain B: 



5 Refinement protocol and experimental data overview

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

Of the 30 calculated structures, 11 were deposited, based on the following criterion: *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
Amber	structure solution	8.0
Amber	refinement	8.0

No chemical shift data was provided.

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:
AAB

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	3.12±0.52	31±11/275 (11.3± 3.9%)	3.97±0.52	76±18/420 (18.0± 4.2%)
2	B	3.25±0.57	40±14/293 (13.5± 4.9%)	3.94±0.53	81±20/451 (18.0± 4.5%)
All	All	3.23	777/6248 (12.4%)	3.99	1725/9581 (18.0%)

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	0.0±0.0	6.2±2.2
2	B	0.0±0.0	6.1±2.2
All	All	0	135

5 of 354 unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
2	B	23	DT	C5-C7	14.26	1.58	1.50	3	5
2	B	19	DT	C5-C7	14.17	1.58	1.50	9	5
1	A	12	DG	P-O5'	13.43	1.73	1.59	4	5
2	B	24	DT	C5-C7	13.10	1.57	1.50	4	4
1	A	8	DA	N3-C4	13.05	1.42	1.34	2	5

5 of 508 unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	20	DG	O4'-C1'-N9	18.44	120.91	108.00	8	4

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	24	DT	C6-C5-C7	-17.76	112.25	122.90	10	8
2	B	15	DC	N3-C2-O2	-17.45	109.68	121.90	10	9
1	A	13	DG	N1-C6-O6	-17.18	109.59	119.90	8	5
2	B	25	DG	C5-C6-N1	17.13	120.07	111.50	1	3

There are no chirality outliers.

5 of 25 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	A	13	DG	Sidechain	10
1	A	11	DG	Sidechain	8
2	B	17	DG	Sidechain	8
2	B	22	DT	Sidechain	8
2	B	26	DG	Sidechain	8

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	5720	3201	3101	-

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 -.

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.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
1	AAB	A	7	1	10,12,13	1.20±0.37	1±1 (10±9%)

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.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
1	AAB	A	7	1	7,16,19	1.59±0.47	1±0 (14±6%)

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	AAB	A	7	1	-	0±0,3,17,18	0±0,1,1,1

5 of 7 unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	7	AAB	C2'-C1'	4.53	1.58	1.52	8	3
1	A	7	AAB	C5'-C4'	3.45	1.62	1.51	9	2
1	A	7	AAB	P-O5'	3.02	1.54	1.62	10	2
1	A	7	AAB	O3'-C3'	2.56	1.48	1.43	7	1
1	A	7	AAB	C2'-C3'	2.44	1.59	1.52	3	2

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	7	AAB	C1'-C2'-C3'	6.08	95.65	103.93	10	10
1	A	7	AAB	O3'-C3'-C2'	2.59	101.65	110.90	6	1

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](#)

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