

wwPDB X-ray Structure Validation Summary Report (i)

May 26, 2020 – 12:11 pm BST

PDB ID : 1D87

Title: CONFORMATIONAL INFLUENCE OF THE RIBOSE 2'-HYDROXYL

GROUP: CRYSTAL STRUCTURES OF DNA-RNA CHIMERIC DUPLEXES

Authors : Egli, M.; Usman, N.; Rich, A.

Deposited on : 1992-08-28

Resolution : 2.25 Å(reported)

This is a wwPDB X-ray 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/XrayValidationReportHelp 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

Xtriage (Phenix) : NOT EXECUTED EDS : NOT EXECUTED

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

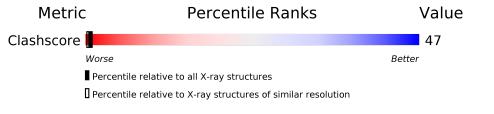
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 2.25 Å.

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	Similar resolution		
Metric	$(\# ext{Entries})$	$(\# ext{Entries}, ext{resolution range}(ext{Å}))$		
Clashscore	141614	1487 (2.26-2.26)		

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	10	10% 90%
1	В	10	10% 90%



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 516 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

• Molecule 1 is DNA/RNA hybrid called DNA/RNA (5'-R(*GP*)-D(*CP*GP*TP*AP*TP* AP*CP*GP*C)-3').

Mol	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
1	Λ	10	Total	С	N	О	Р	0	0	0
1	1 A	10	203	97	38	59	9			
1	D	10	Total	С	N	О	Р	0	0	0
1	Б	10	203	97	38	59	9	U	U	U

• Molecule 2 is water.

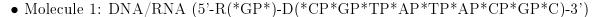
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	61	Total O 61 61	0	0
2	В	49	Total O 49 49	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-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 outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.



Chain A: 10% 90%

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• Molecule 1: DNA/RNA (5'-R(*GP*)-D(*CP*GP*TP*AP*TP*AP*CP*GP*C)-3')

Chain B: 10% 90%



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value
Space group	P 21 21 21
Cell constants	$27.50 \text{\AA} 44.23 \text{Å} 44.61 \text{Å}$
a, b, c, α , β , γ	90.00° 90.00° 90.00°
Resolution (Å)	(Not available) – 2.25
% Data completeness	(Not available) ((Not available)-2.25)
(in resolution range)	(110t available) ((110t available)-2.29)
R_{merge}	(Not available)
R_{sym}	(Not available)
Refinement program	PROLSQ (MODIFIED BY G.J.QUIGLEY), MODIFIED BY G.J.QUIGLEY
R, R_{free}	0.178 , (Not available)
Estimated twinning fraction	No twinning to report.
Total number of atoms	516
Average B, all atoms (Å ²)	22.0



5 Model quality (i)

5.1 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 root-mean-square of all Z scores of the bond lengths (or angles).

Mol Chain		Bo	nd lengths	Bond angles		
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5	
1	A	3.47	$30/227 \ (13.2\%)$	4.74	80/349 (22.9%)	
1	В	3.43	$29/227 \ (12.8\%)$	4.53	72/349 (20.6%)	
All	All	3.45	59/454 (13.0%)	4.63	152/698 (21.8%)	

The worst 5 of 59 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	A	7	DA	C2'-C1'	12.67	1.65	1.52
1	В	19	DG	C2-N2	-11.36	1.23	1.34
1	A	9	DG	C2-N2	-11.33	1.23	1.34
1	В	13	DG	C2-N2	-11.03	1.23	1.34
1	В	11	G	C2-N2	-10.81	1.23	1.34

The worst 5 of 152 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
1	A	7	DA	P-O3'-C3'	26.43	151.41	119.70
1	В	13	DG	O4'-C4'-C3'	-20.86	93.48	106.00
1	A	9	DG	N3-C2-N2	-15.59	108.98	119.90
1	В	12	DC	N3-C4-C5	-13.86	116.36	121.90
1	В	16	DT	O4'-C4'-C3'	-13.83	97.70	106.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Α	203	0	108	11	1
1	В	203	0	108	18	1
2	A	61	0	0	2	5
2	В	49	0	0	2	5
All	All	516	0	216	29	6

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 47.

The worst 5 of 29 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{l} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
1:B:17:DA:H2'	1:B:18:DC:H5"	1.58	0.86
1:B:20:DC:H6	1:B:20:DC:H5'	1.50	0.77
1:A:8:DC:H2'	1:A:9:DG:C8	2.22	0.74
1:B:11:G:H8	1:B:11:G:HO5'	1.38	0.68
1:B:19:DG:C3'	1:B:20:DC:H5"	2.25	0.66

The worst 5 of 6 symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$egin{array}{l} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
2:A:34:HOH:O	2:B:127:HOH:O[1_655]	1.52	0.68
2:A:35:HOH:O	2:B:117:HOH:O[1_655]	1.65	0.55
2:A:24:HOH:O	2:B:52:HOH:O[1_655]	1.75	0.45
2:A:63:HOH:O	2:B:104:HOH:O[4_565]	1.81	0.39
1:A:4:DT:O2	1:B:11:G:N2[2_664]	1.98	0.22

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

There are no protein molecules in this entry.

5.3.2 Protein sidechains (i)

There are no protein molecules in this entry.



5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

