



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 07:01 pm GMT

PDB ID : 220D
Title : INFLUENCE OF COUNTER-IONS ON THE CRYSTAL STRUCTURES OF
DNA DECAMERS: BINDING OF [CO(NH₃)₆]³⁺ AND BA₂⁺ TO A-DNA
Authors : Gao, Y.-G.; Robinson, H.; Van Boom, J.H.; Wang, A.H.-J.
Deposited on : 1995-06-26
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

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:

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

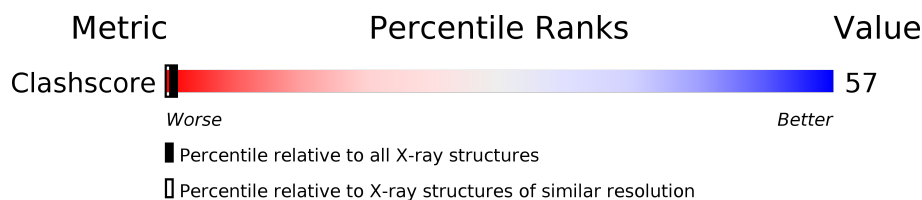
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	7775 (2.00-2.00)

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. 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	10	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 251 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 a DNA chain called DNA (5'-D(*AP*CP*CP*CP*GP*CP*GP*GP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	10	Total	C	N	O	P	0	0	0
			202	96	39	58	9			

- Molecule 2 is BARIUM ION (three-letter code: BA) (formula: Ba).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ba	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	48	Total	O	0	0
			48	48		

3 Residue-property plots

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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

- Molecule 1: DNA (5'-D(*AP*CP*CP*CP*GP*CP*GP*GP*GP*T)-3')

Chain A: 

A1	C2	C3	C4	G5	C6	G7	G8	G9	T10
----	----	----	----	----	----	----	----	----	-----

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	40.12Å 40.12Å 78.22Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	NUCLSQ	Depositor
R, R_{free}	0.206 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	251	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	3.61	42/226 (18.6%)	5.02	71/347 (20.5%)

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	2	DC	P-O5'	-9.63	1.50	1.59
1	A	4	DC	C3'-C2'	9.36	1.63	1.52
1	A	7	DG	C4'-C3'	-8.83	1.43	1.52
1	A	3	DC	C4-C5	-8.73	1.35	1.43
1	A	4	DC	C2'-C1'	-8.66	1.43	1.52
1	A	5	DG	N9-C8	-8.55	1.31	1.37
1	A	5	DG	P-OP1	8.34	1.63	1.49
1	A	6	DC	C2'-C1'	-8.30	1.44	1.52
1	A	4	DC	P-O5'	8.10	1.67	1.59
1	A	6	DC	N3-C4	7.93	1.39	1.33
1	A	3	DC	C4'-O4'	-7.65	1.37	1.45
1	A	1	DA	C6-N1	-7.22	1.30	1.35
1	A	6	DC	C4'-O4'	-7.11	1.38	1.45
1	A	4	DC	O3'-P	-7.03	1.52	1.61
1	A	6	DC	P-O5'	-6.80	1.52	1.59
1	A	6	DC	O4'-C1'	6.66	1.50	1.42
1	A	9	DG	N9-C8	-6.59	1.33	1.37
1	A	1	DA	N7-C5	-6.44	1.35	1.39
1	A	6	DC	C5'-C4'	6.43	1.58	1.51
1	A	7	DG	N9-C8	-6.41	1.33	1.37
1	A	7	DG	C5'-C4'	6.40	1.58	1.51
1	A	8	DG	C4'-C3'	-6.30	1.46	1.52
1	A	8	DG	C5'-C4'	6.22	1.58	1.51
1	A	8	DG	P-OP1	6.09	1.59	1.49
1	A	6	DC	P-OP1	6.06	1.59	1.49
1	A	2	DC	C2-N3	6.00	1.40	1.35

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	4	DC	O4'-C1'	5.88	1.49	1.42
1	A	9	DG	C2'-C1'	-5.85	1.46	1.52
1	A	7	DG	N3-C4	5.78	1.39	1.35
1	A	7	DG	C5-C4	-5.67	1.34	1.38
1	A	4	DC	C4-C5	-5.59	1.38	1.43
1	A	6	DC	C3'-C2'	5.46	1.58	1.52
1	A	8	DG	N9-C8	-5.39	1.34	1.37
1	A	7	DG	N1-C2	-5.39	1.33	1.37
1	A	9	DG	C6-O6	5.35	1.28	1.24
1	A	7	DG	C3'-C2'	5.28	1.58	1.52
1	A	4	DC	C4'-C3'	-5.24	1.47	1.52
1	A	10	DT	C2'-C1'	5.19	1.57	1.52
1	A	10	DT	P-O5'	5.17	1.65	1.59
1	A	6	DC	N1-C2	-5.11	1.35	1.40
1	A	8	DG	C4'-O4'	-5.09	1.40	1.45
1	A	2	DC	P-OP1	-5.06	1.40	1.49

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	5	DG	O5'-P-OP2	25.09	140.81	110.70
1	A	6	DC	O5'-P-OP2	21.79	136.84	110.70
1	A	9	DG	P-O3'-C3'	20.21	143.96	119.70
1	A	1	DA	N1-C6-N6	16.61	128.57	118.60
1	A	5	DG	OP1-P-OP2	-13.76	98.97	119.60
1	A	10	DT	N3-C4-O4	-13.62	111.73	119.90
1	A	7	DG	OP2-P-O3'	13.55	135.01	105.20
1	A	5	DG	N1-C6-O6	-13.17	112.00	119.90
1	A	7	DG	N1-C6-O6	-13.05	112.07	119.90
1	A	10	DT	C2-N3-C4	-13.01	119.40	127.20
1	A	4	DC	OP1-P-OP2	12.87	138.90	119.60
1	A	3	DC	P-O5'-C5'	-12.59	100.76	120.90
1	A	10	DT	OP1-P-OP2	12.43	138.24	119.60
1	A	5	DG	P-O5'-C5'	-11.90	101.86	120.90
1	A	5	DG	C6-N1-C2	-11.70	118.08	125.10
1	A	10	DT	N3-C4-C5	11.55	122.13	115.20
1	A	2	DC	C5-C4-N4	11.12	127.99	120.20
1	A	8	DG	O5'-P-OP2	-11.01	95.80	105.70
1	A	1	DA	C5-C6-N6	-10.83	115.03	123.70
1	A	6	DC	OP1-P-OP2	-10.77	103.44	119.60
1	A	1	DA	N1-C2-N3	-10.11	124.25	129.30
1	A	6	DC	C6-N1-C2	10.03	124.31	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	6	DC	O5'-P-OP1	-9.83	96.85	105.70
1	A	8	DG	N1-C6-O6	-9.79	114.03	119.90
1	A	5	DG	C5-C6-N1	9.37	116.19	111.50
1	A	7	DG	OP1-P-OP2	9.34	133.60	119.60
1	A	5	DG	O5'-P-OP1	-9.15	97.47	105.70
1	A	3	DC	P-O3'-C3'	-9.12	108.75	119.70
1	A	7	DG	O5'-P-OP1	-8.79	97.79	105.70
1	A	8	DG	C5-C6-N1	8.74	115.87	111.50
1	A	6	DC	P-O3'-C3'	-8.73	109.22	119.70
1	A	9	DG	C4-C5-N7	-8.58	107.37	110.80
1	A	2	DC	O4'-C4'-C3'	-8.48	100.91	106.00
1	A	2	DC	N3-C4-N4	-8.20	112.26	118.00
1	A	6	DC	O4'-C4'-C3'	-8.05	101.17	106.00
1	A	6	DC	O4'-C1'-N1	-8.02	102.39	108.00
1	A	3	DC	O3'-P-O5'	-7.95	88.90	104.00
1	A	9	DG	N1-C6-O6	-7.77	115.24	119.90
1	A	7	DG	C5-C6-N1	7.66	115.33	111.50
1	A	9	DG	N3-C2-N2	-7.39	114.72	119.90
1	A	9	DG	P-O5'-C5'	-7.35	109.14	120.90
1	A	1	DA	C6-N1-C2	7.33	123.00	118.60
1	A	7	DG	P-O5'-C5'	-7.30	109.22	120.90
1	A	4	DC	P-O5'-C5'	-7.24	109.31	120.90
1	A	1	DA	OP2-P-O3'	-6.95	89.91	105.20
1	A	2	DC	O4'-C1'-N1	-6.93	103.15	108.00
1	A	1	DA	OP1-P-O3'	6.86	120.30	105.20
1	A	8	DG	O5'-P-OP1	-6.81	99.57	105.70
1	A	9	DG	O4'-C1'-N9	6.69	112.69	108.00
1	A	3	DC	N3-C4-N4	-6.67	113.33	118.00
1	A	7	DG	C5-C6-O6	6.58	132.55	128.60
1	A	9	DG	O5'-P-OP2	6.42	118.41	110.70
1	A	8	DG	C6-C5-N7	6.39	134.24	130.40
1	A	5	DG	N3-C4-C5	-6.12	125.54	128.60
1	A	8	DG	C6-N1-C2	-6.10	121.44	125.10
1	A	10	DT	P-O5'-C5'	-6.05	111.22	120.90
1	A	3	DC	O5'-P-OP2	5.89	117.77	110.70
1	A	10	DT	O4'-C1'-N1	5.82	112.08	108.00
1	A	10	DT	C5-C6-N1	-5.79	120.22	123.70
1	A	9	DG	C6-C5-N7	5.77	133.86	130.40
1	A	9	DG	N9-C4-C5	5.76	107.70	105.40
1	A	10	DT	O5'-P-OP1	-5.58	100.67	105.70
1	A	10	DT	N1-C2-N3	5.55	117.93	114.60
1	A	9	DG	C5-C6-O6	5.49	131.90	128.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2	DC	N1-C2-O2	5.46	122.18	118.90
1	A	5	DG	N1-C2-N3	5.46	127.18	123.90
1	A	7	DG	N3-C2-N2	5.38	123.66	119.90
1	A	2	DC	N3-C4-C5	-5.37	119.75	121.90
1	A	5	DG	C4-C5-N7	-5.36	108.66	110.80
1	A	5	DG	C5-C6-O6	5.26	131.75	128.60
1	A	6	DC	C3'-C2'-C1'	-5.22	96.23	102.50

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	A	202	0	113	18	0
2	A	1	0	0	0	0
3	A	48	0	0	0	4
All	All	251	0	113	18	4

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 57.

All (18) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:DG:H2''	1:A:10:DT:H5'	1.43	0.98
1:A:9:DG:C2'	1:A:10:DT:H5'	2.15	0.75
1:A:5:DG:H5'	1:A:5:DG:H8	1.49	0.75
1:A:3:DC:H2'	1:A:4:DC:C6	2.30	0.67
1:A:8:DG:H2'	1:A:9:DG:H5'	1.79	0.65
1:A:8:DG:H5''	1:A:8:DG:H8	1.63	0.62
1:A:5:DG:H5'	1:A:5:DG:C8	2.35	0.59
1:A:8:DG:H2'	1:A:9:DG:C5'	2.35	0.56
1:A:9:DG:C2'	1:A:10:DT:C5'	2.87	0.52
1:A:10:DT:C6	1:A:10:DT:H3'	2.47	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:DG:H2''	1:A:10:DT:C5'	2.28	0.49
1:A:3:DC:H2'	1:A:4:DC:H6	1.75	0.49
1:A:10:DT:C6	1:A:10:DT:C3'	2.97	0.48
1:A:8:DG:H5''	1:A:8:DG:C8	2.48	0.45
1:A:8:DG:C2'	1:A:9:DG:C5'	2.97	0.43
1:A:8:DG:C2'	1:A:9:DG:H5'	2.47	0.42
1:A:4:DC:H2'	1:A:5:DG:H5'	2.02	0.42
1:A:6:DC:C2'	1:A:7:DG:H5'	2.51	0.40

All (4) 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	Interatomic distance (Å)	Clash overlap (Å)
3:A:14:HOH:O	3:A:14:HOH:O[11_555]	1.36	0.84
3:A:24:HOH:O	3:A:24:HOH:O[12_555]	1.83	0.37
3:A:53:HOH:O	3:A:53:HOH:O[10_665]	1.84	0.36
3:A:26:HOH:O	3:A:26:HOH:O[12_565]	1.93	0.27

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](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.