

wwPDB X-ray Structure Validation Summary Report (i)

Feb 6, 2024 – 03:14 PM EST

PDB ID : 2GB9

Title: d(CGTACG)2 crosslinked bis-acridine complex

Authors: Hopcroft, N.H.; Brogden, A.L.; Searcey, M.; Cardin, C.J.

Deposited on : 2006-03-10

Resolution : 1.70 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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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)

Xtriage (Phenix) : 1.13

EDS : 2.36

buster-report : 1.1.7 (2018)

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

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

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

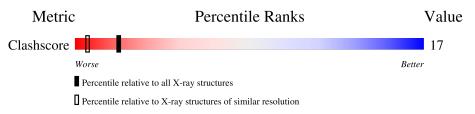
Validation Pipeline (wwPDB-VP) : 2.36

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 1.70 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	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	4695 (1.70-1.70)

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 of 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%

Mol	Chain	Length	Quality of chain			
1	A	6	67%	33%		
1	В	6	100%			



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 397 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.

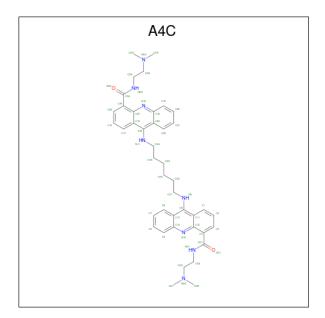
• Molecule 1 is a DNA chain called 5'-D(*CP*GP*TP*AP*CP*G)-3'.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	6	Total	С	N	О	Р	0	0	0
1	A	U	120	58	23	34	5	U		U
1	D	6	Total	С	N	О	Р	0	0	0
1	Б	0	120	58	23	34	5	U		0

• Molecule 2 is STRONTIUM ION (three-letter code: SR) (formula: Sr).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Sr 1 1	0	0
2	В	1	Total Sr 1 1	0	0

• Molecule 3 is 9,9'-(HEXANE-1,6-DIYLDIIMINO)BIS{N-[2-(DIMETHYLAMINO)ETHYL] ACRIDINE-4-CARBOXAMIDE} (three-letter code: A4C) (formula: C₄₂H₅₀N₈O₂).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	В	1	Total 26	_		_	0	0
3	В	1	Total 26	C 21	N 4	O 1	0	0

\bullet Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	48	Total O 48 48	0	0
4	В	55	Total O 55 55	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

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

Chain A: 67% 33%

C1 G2 T3 A4 C5 G6

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

Chain B: 100%

C1 G2 T3 A4 C5 G6



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants	37.21Å 37.21Å 53.48Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 - 1.70	Depositor
rtesolution (A)	30.54 - 1.70	EDS
% Data completeness	95.0 (8.00-1.70)	Depositor
(in resolution range)	100.0 (30.54-1.70)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	13.75 (at 1.70Å)	Xtriage
Refinement program	SHELXL-97	Depositor
D D.	0.213 , 0.254	Depositor
R, R_{free}	0.252 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	15.4	Xtriage
Anisotropy	0.292	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.38, 58.4	EDS
L-test for twinning ²	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	397	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 17.13% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: SR, A4C

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	2.20	7/134 (5.2%)	3.42	$23/205 \ (11.2\%)$	
1	В	2.25	8/134 (6.0%)	3.20	19/205 (9.3%)	
All	All	2.23	15/268 (5.6%)	3.31	42/410 (10.2%)	

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\operatorname{Ideal}(\text{\AA})$
1	A	3	DT	O4'-C1'	7.08	1.50	1.42
1	A	6	DG	C1'-N9	6.38	1.57	1.49
1	В	2	DG	C1'-N9	6.37	1.57	1.49
1	В	6	DG	C1'-N9	6.32	1.57	1.49
1	A	4	DA	O4'-C1'	6.20	1.49	1.42

The worst 5 of 42 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	5	DC	O4'-C1'-N1	17.55	120.28	108.00
1	A	6	DG	O4'-C1'-N9	-16.55	96.42	108.00
1	В	2	DG	O4'-C1'-N9	-14.83	97.62	108.00
1	В	1	DC	C6-N1-C2	12.18	125.17	120.30
1	В	3	DT	N3-C2-O2	11.92	129.45	122.30

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	$\mathbf{H}(\mathbf{model})$	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
1	A	120	0	69	2	0
1	В	120	0	68	0	0
2	A	1	0	0	0	0
2	В	1	0	0	0	0
3	В	52	0	46	7	0
4	A	48	0	0	1	0
4	В	55	0	0	1	0
All	All	397	0	183	8	0

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 17.

The worst 5 of 8 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
3:B:302:A4C:HX21	3:B:302:A4C:H8	1.76	0.67
3:B:302:A4C:HX11	4:B:321:HOH:O	2.00	0.62
3:B:301:A4C:HX21	3:B:301:A4C:H8	1.87	0.56
1:A:5:DC:H2"	3:B:301:A4C:H7	1.96	0.47
1:A:4:DA:H4'	4:A:245:HOH:O	2.15	0.46

There are no symmetry-related clashes.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or 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 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| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mal	Mol Type Chain Res		Dag	Link	Bond lengths			Bond angles		
MIOI	Type	Chain	Res	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	A4C	В	301	3	28,28,57	1.10	2 (7%)	37,38,78	1.42	6 (16%)
3	A4C	В	302	3	28,28,57	1.04	1 (3%)	37,38,78	1.35	5 (13%)

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.

N	Лol	Type	Chain	Res	Link	Chirals	Torsions	Rings
	3	A4C	В	301	3	-	7/14/14/31	0/3/3/6
	3	A4C	В	302	3	-	7/14/14/31	0/3/3/6

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
3	В	301	A4C	C4-C12	-2.73	1.39	1.43
3	В	302	A4C	C4-C12	-2.47	1.39	1.43
3	В	301	A4C	C13-C14	-2.00	1.39	1.42

The worst 5 of 11 bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
3	В	301	A4C	C12-N10-C14	4.43	122.07	117.02
3	В	302	A4C	C12-N10-C14	3.97	121.54	117.02
3	В	301	A4C	C11-C12-N10	-2.80	119.44	123.31
3	В	302	A4C	C1-C11-C9	-2.58	118.92	123.40
3	В	302	A4C	C13-C14-N10	-2.57	120.94	123.35

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	В	301	A4C	C3-C4-CD1-OD1
3	В	301	A4C	C12-C4-CD1-OD1
3	В	301	A4C	C3-C4-CD1-ND1
3	В	302	A4C	C3-C4-CD1-OD1
3	В	302	A4C	C12-C4-CD1-OD1

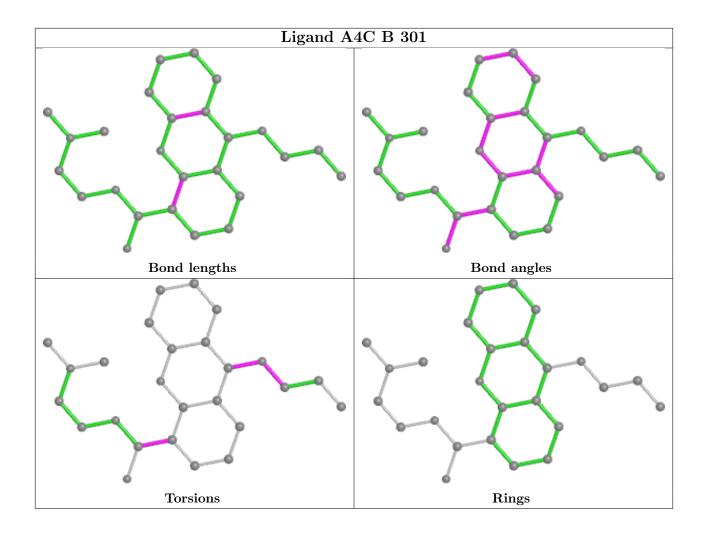
There are no ring outliers.

2 monomers are involved in 7 short contacts:

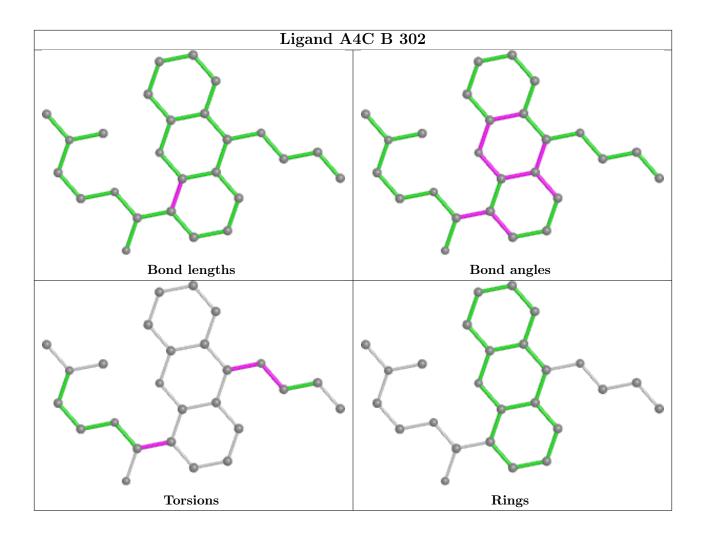
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	В	301	A4C	3	0
3	В	302	A4C	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









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)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

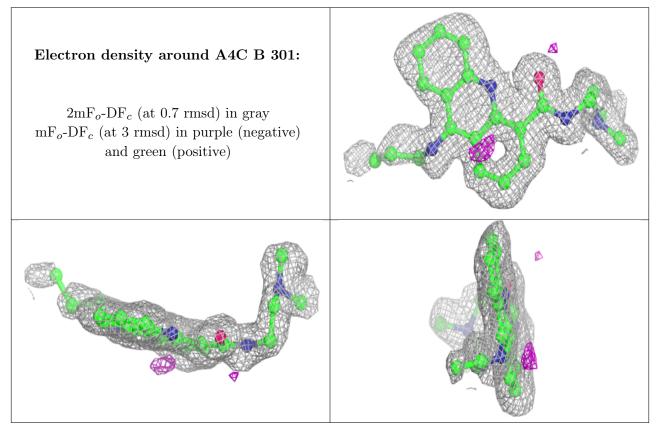
6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

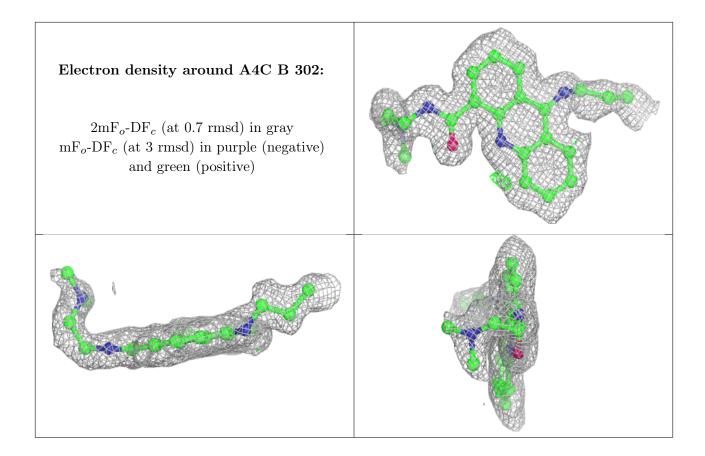
6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.







6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

