



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 11:04 am BST

PDB ID : 272D
Title : PARALLEL AND ANTIPARALLEL (G.GC)₂ TRIPLE HELIX FRAGMENTS IN A CRYSTAL STRUCTURE
Authors : Vlieghe, D.; Van Meervelt, L.; Dautant, A.; Gallois, B.; Precigoux, G.; Kennard, O.
Deposited on : 1996-07-09
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

<https://www.wwpdb.org/validation/2017/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) : 1.13
EDS : 2.11
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) : 2.11

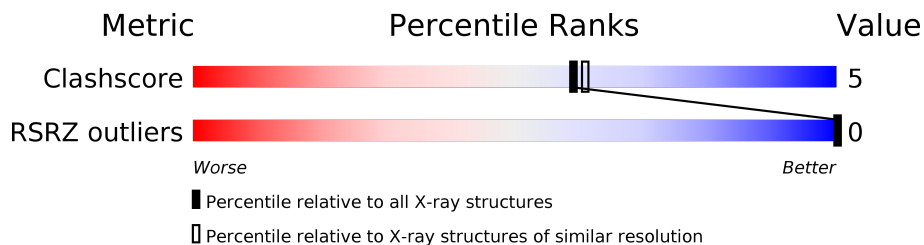
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

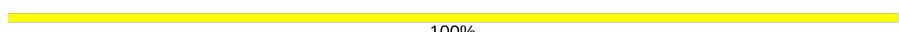
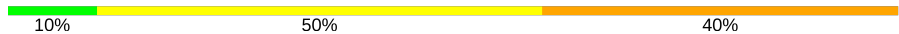
The reported resolution of this entry is 2.00 Å.

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	141614	9178 (2.00-2.00)
RSRZ outliers	127900	7900 (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 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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	A	10	 100%
1	B	10	 10% 50% 40%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 454 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(*GP*GP*CP*CP*AP*AP*TP*TP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	10	Total	C	N	O	P	0	0	0
			205	98	40	58	9			
1	B	10	Total	C	N	O	P	0	0	0
			205	98	40	58	9			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	22	Total	O	0	0
			22	22		
2	B	22	Total	O	0	0
			22	22		

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.

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

Chain A:  100%

G1
G2
C3
C4
A5
A6
T7
T8
G9
G10

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

Chain B:  10% 50% 40%

G11
G12
C13
C14
A15
A16
T17
T18
G19
G20

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	26.25Å 36.82Å 53.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00 13.65 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.00) 93.5 (13.65-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.79 (at 1.99Å)	Xtrriage
Refinement program	SHELXL-93	Depositor
R, R_{free}	0.209 , (Not available) 0.197 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	32.9	Xtrriage
Anisotropy	0.244	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 66.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	454	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

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	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.73	3/230 (1.3%)	2.89	29/354 (8.2%)
1	B	1.59	0/230	2.73	24/354 (6.8%)
All	All	1.66	3/460 (0.7%)	2.81	53/708 (7.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	2	DG	C4'-O4'	-5.78	1.39	1.45
1	A	3	DC	C4'-O4'	-5.71	1.39	1.45
1	A	1	DG	C4'-O4'	-5.19	1.39	1.45

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	DT	O4'-C1'-N1	-13.55	98.52	108.00
1	B	17	DT	O4'-C1'-N1	-13.40	98.62	108.00
1	A	4	DC	O4'-C1'-N1	13.04	117.13	108.00
1	A	6	DA	O4'-C1'-N9	-12.07	99.55	108.00
1	B	16	DA	O4'-C1'-N9	-11.37	100.04	108.00
1	A	7	DT	O4'-C1'-N1	-11.24	100.13	108.00
1	B	18	DT	O4'-C1'-N1	-10.71	100.50	108.00
1	B	19	DG	O4'-C1'-N9	-10.13	100.91	108.00
1	A	5	DA	O4'-C1'-N9	-9.56	101.31	108.00
1	B	15	DA	O4'-C1'-N9	-9.51	101.35	108.00
1	B	13	DC	C2-N1-C1'	-9.46	108.40	118.80
1	A	3	DC	C2-N1-C1'	-8.79	109.13	118.80
1	A	1	DG	O4'-C1'-N9	-8.67	101.93	108.00
1	B	11	DG	O4'-C1'-N9	-8.14	102.30	108.00
1	B	14	DC	C2-N1-C1'	-7.76	110.26	118.80
1	B	11	DG	N1-C6-O6	7.63	124.48	119.90
1	B	13	DC	C6-N1-C1'	7.59	129.91	120.80
1	A	2	DG	O4'-C1'-N9	-7.41	102.81	108.00
1	B	14	DC	C6-N1-C1'	7.30	129.56	120.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	DT	N1-C1'-C2'	6.92	125.75	112.60
1	B	11	DG	C5-C6-O6	-6.91	124.45	128.60
1	B	17	DT	N1-C1'-C2'	6.80	125.53	112.60
1	A	4	DC	C2-N1-C1'	-6.63	111.50	118.80
1	A	3	DC	C6-N1-C2	6.50	122.90	120.30
1	A	2	DG	C8-N9-C4	6.31	108.93	106.40
1	A	7	DT	N1-C1'-C2'	6.29	124.55	112.60
1	B	18	DT	O4'-C1'-C2'	-6.23	100.92	105.90
1	B	15	DA	N1-C6-N6	6.18	122.31	118.60
1	A	2	DG	N9-C1'-C2'	6.08	124.15	112.60
1	A	3	DC	C6-N1-C1'	5.97	127.97	120.80
1	B	12	DG	C5-C6-O6	-5.97	125.02	128.60
1	A	4	DC	C6-N1-C1'	5.97	127.96	120.80
1	B	18	DT	N1-C1'-C2'	5.96	123.93	112.60
1	A	4	DC	N1-C2-O2	5.84	122.40	118.90
1	A	7	DT	O4'-C1'-C2'	-5.81	101.25	105.90
1	A	4	DC	O4'-C1'-C2'	-5.80	101.26	105.90
1	A	5	DA	P-O3'-C3'	5.65	126.48	119.70
1	B	14	DC	OP1-P-OP2	-5.63	111.16	119.60
1	B	16	DA	OP1-P-OP2	-5.55	111.28	119.60
1	A	5	DA	O5'-P-OP2	5.47	117.27	110.70
1	B	18	DT	OP1-P-OP2	-5.38	111.54	119.60
1	B	12	DG	OP1-P-OP2	-5.30	111.65	119.60
1	B	14	DC	O4'-C1'-N1	5.23	111.66	108.00
1	A	10	DG	O4'-C1'-N9	-5.18	104.38	108.00
1	A	2	DG	N3-C2-N2	5.17	123.52	119.90
1	A	9	DG	O4'-C1'-N9	-5.14	104.40	108.00
1	A	6	DA	OP1-P-OP2	-5.12	111.92	119.60
1	A	4	DC	C3'-C2'-C1'	-5.09	96.39	102.50
1	A	5	DA	N9-C1'-C2'	5.09	122.27	112.60
1	A	3	DC	O5'-P-OP1	5.09	116.81	110.70
1	B	12	DG	O4'-C1'-C2'	-5.07	101.85	105.90
1	A	4	DC	P-O3'-C3'	5.05	125.76	119.70
1	B	15	DA	OP1-P-OP2	-5.02	112.07	119.60

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	205	0	114	0	0
1	B	205	0	114	3	0
2	A	22	0	0	0	0
2	B	22	0	0	0	0
All	All	454	0	228	3	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 5.

All (3) 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:B:15:DA:H2''	1:B:16:DA:OP2	1.92	0.68
1:B:13:DC:H2'	1:B:14:DC:C6	2.44	0.52
1:B:15:DA:C2'	1:B:16:DA:OP2	2.58	0.43

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 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](#)

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q < 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	10/10 (100%)	-0.27	0 100 100	23, 29, 33, 33	0
1	B	10/10 (100%)	-0.38	0 100 100	24, 29, 34, 36	0
All	All	20/20 (100%)	-0.33	0 100 100	23, 29, 34, 36	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.