



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 06:36 am BST

PDB ID : 1QBP
Title : CRYSTAL STRUCTURE OF A BROMINATED RNA HELIX WITH FOUR
MISMATCHED BASE PAIRS
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Deposited on : 1999-04-26
Resolution : 2.10 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

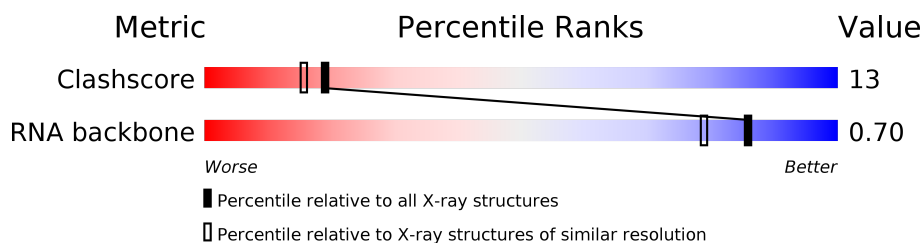
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5710 (2.10-2.10)
RNA backbone	3102	1000 (2.54-1.66)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	15	
1	B	15	
1	C	15	
1	D	15	
1	E	15	
1	F	15	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1910 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 RNA chain called 5'-R(*UP*GP*(CBV)P*(CBV)P*AP*GP*UP*UP*CP*GP*CP*UP*GP*GP*C)-3'.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	15	Total 300	Br 2	C 132	N 51	O 101	P 14	0	0	1
1	B	15	Total 300	Br 2	C 132	N 51	O 101	P 14	0	0	1
1	C	15	Total 300	Br 2	C 132	N 51	O 101	P 14	0	0	1
1	D	15	Total 300	Br 2	C 132	N 51	O 101	P 14	0	0	1
1	E	15	Total 300	Br 2	C 132	N 51	O 101	P 14	0	0	1
1	F	15	Total 300	Br 2	C 132	N 51	O 101	P 14	0	0	1

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	25	Total 25	O 25	0	0
2	B	22	Total 22	O 22	0	0
2	C	32	Total 32	O 32	0	0
2	D	23	Total 23	O 23	0	0
2	E	6	Total 6	O 6	0	0
2	F	2	Total 2	O 2	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.

- Molecule 1: 5'-R(*UP*GP*(CBV)P*(CBV)P*AP*GP*UP*UP*CP*GP*CP*UP*GP*GP*C)-3',

Chain A: 



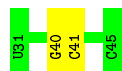
- Molecule 1: 5'-R(*UP*GP*(CBV)P*(CBV)P*AP*GP*UP*UP*CP*GP*CP*UP*GP*GP*C)-3',

Chain B: 



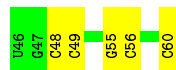
- Molecule 1: 5'-R(*UP*GP*(CBV)P*(CBV)P*AP*GP*UP*UP*CP*GP*CP*UP*GP*GP*C)-3',

Chain C: 



- Molecule 1: 5'-R(*UP*GP*(CBV)P*(CBV)P*AP*GP*UP*UP*CP*GP*CP*UP*GP*GP*C)-3',

Chain D: 



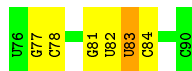
- Molecule 1: 5'-R(*UP*GP*(CBV)P*(CBV)P*AP*GP*UP*UP*CP*GP*CP*UP*GP*GP*C)-3',

Chain E: 



- Molecule 1: 5'-R(*UP*GP*(CBV)P*(CBV)P*AP*GP*UP*UP*CP*GP*CP*UP*GP*GP*C)-3',

Chain F: 60% 33% 7%



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	104.45Å 24.27Å 104.45Å 90.00° 118.76° 90.00°	Depositor
Resolution (Å)	999.00 – 2.10	Depositor
% Data completeness (in resolution range)	97.0 (999.00-2.10)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.217 , 0.253	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1910	wwPDB-VP
Average B, all atoms (Å ²)	19.0	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: CBV

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	1.05	0/286	0.85	0/443
1	B	0.96	1/286 (0.3%)	0.81	0/443
1	C	0.93	0/286	0.82	0/443
1	D	1.12	0/286	0.88	0/443
1	E	0.74	0/286	0.79	0/443
1	F	0.71	0/286	0.80	0/443
All	All	0.93	1/1716 (0.1%)	0.83	0/2658

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	24	C	N1-C2	-5.30	1.34	1.40

There are no bond angle outliers.

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	A	300	0	150	1	0
1	B	300	0	150	4	0
1	C	300	0	150	1	0
1	D	300	0	150	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	300	0	150	18	0
1	F	300	0	150	9	0
2	A	25	0	0	0	0
2	B	22	0	0	1	0
2	C	32	0	0	0	0
2	D	23	0	0	0	0
2	E	6	0	0	0	0
2	F	2	0	0	0	0
All	All	1910	0	900	35	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 13.

All (35) 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:E:69:C:C3'	1:E:70:G:H5''	1.81	1.09
1:E:69:C:C2'	1:E:70:G:H5''	1.91	1.01
1:E:69:C:H2'	1:E:70:G:H5''	1.49	0.95
1:E:69:C:H3'	1:E:70:G:H5''	1.51	0.92
1:E:70:G:O2'	1:E:71:C:H5'	1.71	0.89
1:E:69:C:H3'	1:E:70:G:C5'	2.08	0.84
1:F:77:G:H2'	1:F:78:CBV:H6	1.67	0.76
1:F:82:U:H5'	1:F:83:U:OP1	1.90	0.71
1:E:69:C:C3'	1:E:70:G:C5'	2.62	0.70
1:E:70:G:H2'	1:E:71:C:O4'	1.92	0.70
1:E:70:G:H5'	1:E:70:G:H8	1.62	0.65
1:F:83:U:O2	1:F:83:U:H2'	1.97	0.63
1:E:70:G:C2'	1:E:71:C:H5'	2.32	0.60
1:B:27:U:O2'	1:B:28:G:H5'	2.03	0.58
1:F:81:G:H2'	1:F:82:U:O4'	2.05	0.57
1:E:73:G:H2'	1:E:74:G:O4'	2.08	0.54
1:F:83:U:H5'	1:F:84:C:OP2	2.08	0.53
1:B:27:U:C2'	1:B:28:G:H5'	2.39	0.52
1:F:82:U:H2'	1:F:83:U:C6	2.47	0.50
1:E:69:C:H2'	1:E:70:G:C5'	2.33	0.48
1:B:23:U:H1'	2:B:7109:HOH:O	2.14	0.47
1:E:70:G:C6	1:E:71:C:C4	3.03	0.47
1:B:23:U:H2'	1:B:24:C:C6	2.50	0.46
1:F:77:G:H2'	1:F:78:CBV:C6	2.40	0.46
1:F:83:U:C2'	1:F:83:U:O2	2.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:83:U:C5'	1:F:84:C:OP2	2.64	0.46
1:D:60:C:C4	1:E:75:C:C4	3.05	0.45
1:E:70:G:H2'	1:E:71:C:C6	2.52	0.44
1:D:48:CBV:H2'	1:D:49:CBV:H6	2.00	0.43
1:E:73:G:O2'	1:E:74:G:H5'	2.19	0.42
1:E:70:G:C2'	1:E:71:C:C5'	2.98	0.42
1:E:73:G:C2'	1:E:74:G:H5'	2.50	0.41
1:D:55:G:C5	1:D:56:C:C5	3.09	0.41
1:C:40:G:H2'	1:C:41:C:O4'	2.22	0.41
1:A:6:G:H2'	1:A:7:U:O4'	2.21	0.40

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](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	11/15 (73%)	0	0
1	B	11/15 (73%)	0	0
1	C	11/15 (73%)	0	0
1	D	11/15 (73%)	0	0
1	E	11/15 (73%)	1 (9%)	0
1	F	11/15 (73%)	1 (9%)	0
All	All	66/90 (73%)	2 (3%)	0

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	E	70	G
1	F	83	U

There are no RNA pucker outliers to report.

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

12 non-standard protein/DNA/RNA residues are modelled in this entry.

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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
1	CBV	E	64	1	15,22,23	1.38	2 (13%)	18,32,35	1.22	3 (16%)
1	CBV	D	49	1	15,22,23	1.08	1 (6%)	18,32,35	1.41	3 (16%)
1	CBV	A	4	1	15,22,23	1.14	1 (6%)	18,32,35	1.21	2 (11%)
1	CBV	B	18	1	15,22,23	1.07	1 (6%)	18,32,35	1.68	4 (22%)
1	CBV	C	34	1	15,22,23	1.04	1 (6%)	18,32,35	1.54	3 (16%)
1	CBV	F	78	1	15,22,23	1.13	1 (6%)	18,32,35	1.50	4 (22%)
1	CBV	A	3	1	15,22,23	1.07	1 (6%)	18,32,35	1.34	3 (16%)
1	CBV	C	33	1	15,22,23	1.14	1 (6%)	18,32,35	1.37	3 (16%)
1	CBV	E	63	1	15,22,23	0.86	1 (6%)	18,32,35	1.45	3 (16%)
1	CBV	F	79	1	15,22,23	0.98	1 (6%)	18,32,35	1.33	3 (16%)
1	CBV	B	19	1	15,22,23	1.00	1 (6%)	18,32,35	1.56	4 (22%)
1	CBV	D	48	1	15,22,23	1.06	1 (6%)	18,32,35	1.52	3 (16%)

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	CBV	E	64	1	-	0/5/25/26	0/2/2/2
1	CBV	D	49	1	-	0/5/25/26	0/2/2/2
1	CBV	A	4	1	-	0/5/25/26	0/2/2/2
1	CBV	B	18	1	-	0/5/25/26	0/2/2/2
1	CBV	C	34	1	-	0/5/25/26	0/2/2/2
1	CBV	F	78	1	-	0/5/25/26	0/2/2/2
1	CBV	A	3	1	-	0/5/25/26	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CBV	C	33	1	-	0/5/25/26	0/2/2/2
1	CBV	E	63	1	-	0/5/25/26	0/2/2/2
1	CBV	F	79	1	-	0/5/25/26	0/2/2/2
1	CBV	B	19	1	-	0/5/25/26	0/2/2/2
1	CBV	D	48	1	-	0/5/25/26	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	48	CBV	C6-C5	-3.43	1.32	1.39
1	E	64	CBV	C6-C5	-3.43	1.32	1.39
1	F	78	CBV	C6-C5	-3.17	1.33	1.39
1	A	3	CBV	C6-C5	-2.89	1.33	1.39
1	C	34	CBV	C6-C5	-2.84	1.33	1.39
1	A	4	CBV	C6-C5	-2.83	1.33	1.39
1	B	18	CBV	C6-C5	-2.55	1.34	1.39
1	C	33	CBV	C6-C5	-2.55	1.34	1.39
1	E	63	CBV	C6-C5	-2.53	1.34	1.39
1	B	19	CBV	C6-C5	-2.51	1.34	1.39
1	D	49	CBV	C6-C5	-2.41	1.34	1.39
1	E	64	CBV	C2-N3	-2.35	1.33	1.38
1	F	79	CBV	C4-C5	2.32	1.45	1.41

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	48	CBV	C5-C4-N4	-4.36	119.33	122.94
1	C	34	CBV	C5-C4-N4	-4.01	119.62	122.94
1	E	63	CBV	C5-C4-N4	-3.88	119.73	122.94
1	B	18	CBV	BR-C5-C6	3.81	126.09	117.34
1	B	18	CBV	BR-C5-C4	-3.80	116.34	120.15
1	F	78	CBV	C5-C4-N4	-3.65	119.92	122.94
1	D	49	CBV	C5-C4-N4	-3.50	120.04	122.94
1	B	19	CBV	C2-N3-C4	3.45	120.18	116.02
1	D	49	CBV	C2-N3-C4	3.42	120.15	116.02
1	A	3	CBV	C5-C4-N4	-3.36	120.16	122.94
1	B	19	CBV	C5-C4-N4	-3.35	120.17	122.94
1	A	4	CBV	C2-N3-C4	3.35	120.06	116.02
1	C	33	CBV	BR-C5-C4	3.28	123.43	120.15
1	F	79	CBV	BR-C5-C4	3.22	123.37	120.15
1	B	18	CBV	C2-N3-C4	3.18	119.85	116.02
1	C	34	CBV	BR-C5-C4	3.16	123.31	120.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	79	CBV	C2-N3-C4	3.12	119.78	116.02
1	E	63	CBV	C2-N3-C4	3.10	119.76	116.02
1	B	19	CBV	BR-C5-C6	3.07	124.40	117.34
1	F	78	CBV	C2-N3-C4	3.07	119.72	116.02
1	C	33	CBV	C2-N3-C4	3.06	119.71	116.02
1	B	18	CBV	C5-C4-N4	-3.04	120.43	122.94
1	E	64	CBV	C5-C4-N4	-2.97	120.48	122.94
1	E	64	CBV	C2-N3-C4	2.90	119.52	116.02
1	C	33	CBV	C5-C4-N4	-2.89	120.55	122.94
1	A	3	CBV	C2-N3-C4	2.87	119.48	116.02
1	F	78	CBV	BR-C5-C6	2.87	123.94	117.34
1	D	48	CBV	C2-N3-C4	2.85	119.46	116.02
1	C	34	CBV	C2-N3-C4	2.83	119.44	116.02
1	A	4	CBV	C5-C4-N4	-2.80	120.62	122.94
1	D	48	CBV	BR-C5-C6	2.58	123.27	117.34
1	E	63	CBV	BR-C5-C6	2.52	123.15	117.34
1	F	79	CBV	C5-C4-N4	-2.41	120.94	122.94
1	E	64	CBV	BR-C5-C6	2.38	122.82	117.34
1	D	49	CBV	BR-C5-C6	2.33	122.69	117.34
1	A	3	CBV	BR-C5-C4	2.20	122.35	120.15
1	B	19	CBV	BR-C5-C4	-2.11	118.03	120.15
1	F	78	CBV	BR-C5-C4	-2.05	118.09	120.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	49	CBV	1	0
1	F	78	CBV	2	0
1	D	48	CBV	1	0

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

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.