

Full wwPDB X-ray Structure Validation Report (i)

May 14, 2020 – 05:31 am BST

PDB ID : 3GOT

> Title : Guanine riboswitch C74U mutant bound to 2-fluoroadenine.

Authors Gilbert, S.D.; Reyes, F.E.; Batey, R.T.

2009-03-20 Deposited on

1.95 Å(reported) Resolution

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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity 4.02b-467

> 1.8.5 (274361), CSD as541be (2020) Mogul

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)

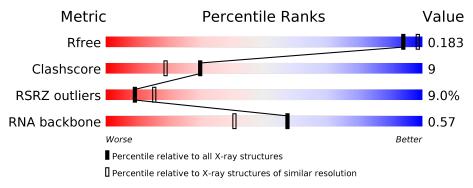
Validation Pipeline (wwPDB-VP) 2.11

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.95 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	$(\# ext{Entries})$	$(\# ext{Entries}, ext{resolution range}(ext{Å}))$
R_{free}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)
RNA backbone	3102	1124 (2.50-1.42)

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% 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	
			9%		
1	A	67	36%	52%	10% •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

M	Iol	Type	Chain	${f Res}$	Chirality	Geometry	Clashes	Electron density
4	4	NCO	A	106	-	-	-	X
4	4	NCO	A	111	-	-	-	X



2 Entry composition (i)

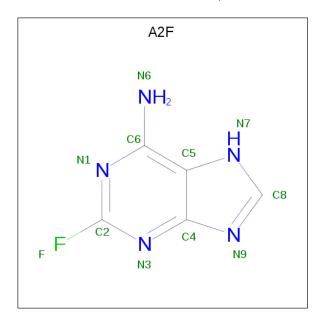
There are 5 unique types of molecules in this entry. The entry contains 1716 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Guanine riboswitch.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	67	Total 1422	C 637	N 252	O 467	P 66	0	0	0

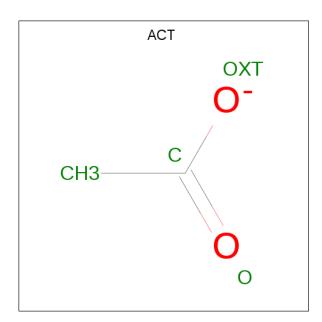
• Molecule 2 is 2-fluoroadenine (three-letter code: A2F) (formula: C₅H₄FN₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 11	C 5	F 1	N 5	0	0

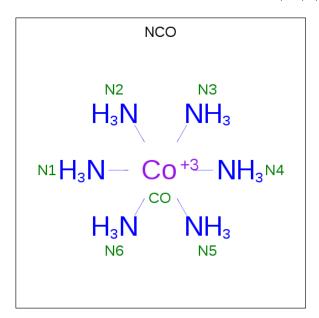
• Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
3	Δ	1	Total	С	О	0	0
	11	1	4	2	2	U	U

 \bullet Molecule 4 is COBALT HEXAMMINE(III) (three-letter code: NCO) (formula: $\mathrm{CoH_{18}N_6}).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0



Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0
4	A	1	Total Co N 7 1 6	0	0

• Molecule 5 is water.

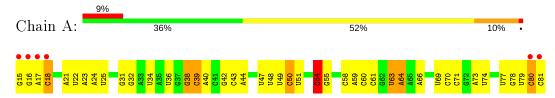
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	202	Total O 202 202	0	0



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 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: Guanine riboswitch





4 Data and refinement statistics (i)

Property	Value	Source	
Space group	C 1 2 1	Depositor	
Cell constants	$132.46 { $	Depositor	
a, b, c, α , β , γ	90.00° 90.10° 90.00°	Depositor	
Resolution (Å)	19.98 - 1.95	Depositor	
Resolution (A)	19.98 - 1.95	EDS	
% Data completeness	95.2 (19.98-1.95)	Depositor	
(in resolution range)	95.4 (19.98-1.95)	EDS	
R_{merge}	(Not available)	Depositor	
R_{sym}	0.06	Depositor	
$< I/\sigma(I) > 1$	$2.68 \; (at \; 1.94 \text{Å})$	Xtriage	
Refinement program	CNS	Depositor	
P. P.	0.218 , 0.268	Depositor	
R, R_{free}	0.155 , 0.183	DCC	
R_{free} test set	134 reflections (0.98%)	wwPDB-VP	
Wilson B-factor (Å ²)	37.1	Xtriage	
Anisotropy	0.102	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.27, 45.2	EDS	
L-test for twinning ²	$< L >=0.52, < L^2>=0.36$	Xtriage	
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage	
F_o, F_c correlation	0.96	EDS	
Total number of atoms	1716	wwPDB-VP	
Average B, all atoms (Å ²)	39.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 9.27% 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: ACT, A2F, NCO

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	Boı	nd lengths	Bond angles		
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	1.16	7/1589 (0.4%)	1.09	7/2474 (0.3%)	

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(\operatorname{\AA})$
1	A	36	U	N1-C2	6.27	1.44	1.38
1	A	66	A	C6-N6	5.91	1.38	1.33
1	A	69	U	C4-O4	5.88	1.28	1.23
1	A	40	A	C6-N6	5.81	1.38	1.33
1	A	71	С	N1-C2	5.65	1.45	1.40
1	A	39	С	C4-N4	5.64	1.39	1.33
1	A	58	С	C2-O2	5.63	1.29	1.24

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\mathbf{Ideal}(^o)$
1	A	39	С	O5'-P-OP2	-13.54	93.52	105.70
1	A	54	С	C5'-C4'-C3'	9.41	131.06	116.00
1	A	39	С	C5'-C4'-C3'	7.91	128.65	116.00
1	A	25	U	O5'-P-OP2	-6.05	100.25	105.70
1	A	63	U	C2'-C3'-O3'	5.89	123.12	113.70
1	A	39	С	O5'-P-OP1	5.47	117.26	110.70



Continued from previous page...

Mol	Chain	Res	Type	${f Atoms}$	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	54	С	C4'-C3'-C2'	-5.15	97.45	102.60

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	34	U	Sidechain
1	A	38	G	Sidechain
1	A	55	G	Sidechain
1	A	64	A	Sidechain
1	A	70	С	Sidechain

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	1422	0	722	21	0
2	A	11	0	4	0	0
3	A	4	0	3	1	0
4	A	77	0	0	0	0
5	A	202	0	0	3	0
All	All	1716	0	729	21	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 9.

All (21) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} & ext{Interatomic} \ & ext{distance} \ & ext{(Å)} \end{aligned}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:A:39:C:H5"	5:A:451:HOH:O	2.00	0.61
1:A:42:G:O2'	1:A:43:C:H5'	2.02	0.59
1:A:23:A:H3'	1:A:24:A:H5'	1.88	0.56
1:A:73:A:H2'	1:A:74:U:O4'	2.06	0.56
1:A:38:G:O2'	3:A:96:ACT:H1	2.09	0.52
1:A:15:G:N3	1:A:15:G:H2'	2.25	0.52



Continued from previous page...

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	$oxed{ ext{distance (Å)} }$	overlap (Å)
1:A:21:A:H2'	1:A:22:U:H2'	1.93	0.50
1:A:17:A:H2'	1:A:18:C:O4'	2.12	0.50
1:A:44:A:H5"	1:A:44:A:C8	2.47	0.49
1:A:32:G:H1'	5:A:353:HOH:O	2.15	0.46
1:A:42:G:HO2'	1:A:43:C:H5'	1.80	0.46
1:A:44:A:H5"	1:A:44:A:H8	1.81	0.45
1:A:79:U:H2'	1:A:80:C:C6	2.54	0.43
1:A:31:G:C2'	1:A:32:G:H5'	2.50	0.42
1:A:54:C:H5"	5:A:351:HOH:O	2.19	0.42
1:A:21:A:O3'	1:A:47:U:H4'	2.20	0.42
1:A:77:U:H2'	1:A:78:G:O4'	2.20	0.41
1:A:50:C:H2'	1:A:51:U:O4'	2.21	0.41
1:A:60:C:H2'	1:A:61:C:C6	2.56	0.40
1:A:59:A:O2'	1:A:60:C:H5'	2.21	0.40
1:A:31:G:O2'	1:A:32:G:H5'	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	66/67~(98%)	9 (13%)	3 (4%)

All (9) RNA backbone outliers are listed below:

Mol	Chain	${ m Res}$	\mathbf{Type}
1	A	16	G
1	A	18	С
1	A	48	U



Continued from previous page...

Mol	Chain	Res	Type
1	A	49	U
1	A	50	С
1	A	54	С
1	A	64	A
1	A	80	С
1	A	81	С

All (3) RNA pucker outliers are listed below:

Mol	Chain	${f Res}$	Type
1	A	48	U
1	A	54	С
1	A	63	U

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)

13 ligands 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	Res Link	B	Bond lengths			Bond angles		
$oxed{f Mol\ Type}$	Chain	Counts			RMSZ	# Z > 2	Counts	RMSZ	# Z > 2		
4	NCO	A	104	-	6,6,6	1.03	1 (16%)	-			
4	NCO	A	107	-	6,6,6	0.25	0	-			
3	ACT	A	96	_	1,3,3	4.97	1 (100%)	0,3,3	0.00	-	
4	NCO	A	101	_	6,6,6	1.32	0	-			
4	NCO	A	108	_	6,6,6	0.08	0	-			



Mol	Trino	Chain	Res	Link	В	ond leng	${ m gths}$	Bond angles	
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	$\mid \text{RMSZ} \mid \# Z > 2$
4	NCO	A	106	-	6,6,6	0.29	0	-	
4	NCO	A	112	-	6,6,6	0.12	0	-	
4	NCO	A	102	_	6,6,6	0.53	0	_	
4	NCO	A	109	_	6,6,6	0.35	0	_	
4	NCO	A	111	-	6,6,6	0.32	0	-	
4	NCO	A	103	_	6,6,6	0.54	0	_	
2	A2F	A	91	-	10,12,12	1.70	2 (20%)	8,17,17	3.93 6 (75%)
4	NCO	A	105	_	6,6,6	0.24	0	-	

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.

\mathbf{Mol}	\mathbf{Type}	Chain	${f Res}$	Link	Chirals	Torsions	Rings
2	A2F	A	91	_	-	_	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	${f Observed(\AA)}$	$\mathbf{Ideal}(\mathbf{\AA})$
3	A	96	ACT	СН3-С	4.97	1.55	1.48
2	A	91	A2F	C2-N1	3.31	1.34	1.31
2	A	91	A2F	C2-N3	3.31	1.34	1.31
4	A	104	NCO	CO-N4	2.22	2.04	1.96

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
2	A	91	A2F	N3-C2-N1	-6.51	121.27	130.51
2	A	91	A2F	F-C2-N1	5.62	120.06	114.73
2	A	91	A2F	C2-N3-C4	4.47	120.47	115.03
2	A	91	A2F	F-C2-N3	4.15	118.67	114.73
2	A	91	A2F	C5-C6-N1	-2.49	119.37	121.01
2	A	91	A2F	C4-C5-N7	-2.16	107.15	109.40

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	96	ACT	1	0

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, 95^{th} 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	67/67 (100%)	0.73	6 (8%) 9 15	22, 33, 74, 93	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	81	С	8.6
1	A	15	G	4.5
1	A	16	G	4.2
1	A	17	A	3.4
1	A	80	С	3.2
1	A	18	С	2.9

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)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B\text{-}factors}({f \AA}^2)$	Q < 0.9
4	NCO	A	106	7/7	0.15	0.94	121,121,121,122	0



Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$ extbf{B-factors}(extbf{A}^2)$	Q<0.9
4	NCO	A	111	7/7	0.48	0.66	103,103,103,104	0
4	NCO	A	107	7/7	0.79	0.31	81,81,81,82	0
2	A2F	A	91	11/11	0.89	0.29	31,34,35,36	0
4	NCO	A	108	7/7	0.90	0.20	67,67,67,67	0
4	NCO	A	109	7/7	0.91	0.33	66,66,67,67	0
4	NCO	A	105	7/7	0.92	0.18	65,65,66,66	0
4	NCO	A	102	7/7	0.93	0.26	$42,\!43,\!44,\!45$	0
4	NCO	A	103	7/7	0.94	0.20	$31,\!34,\!36,\!37$	0
3	ACT	A	96	4/4	0.95	0.31	19,20,22,22	0
4	NCO	A	112	7/7	0.95	0.21	36,36,37,38	7
4	NCO	A	101	7/7	0.98	0.14	23,26,28,28	0
4	NCO	A	104	7/7	0.98	0.13	24,27,29,30	0

6.5 Other polymers (i)

There are no such residues in this entry.

