

# Full wwPDB X-ray Structure Validation Report (i)

Jan 6, 2024 – 07:31 pm GMT

PDB ID : 6H3V

Title: Bunyamwera Virus Glycoprotein Gc Head Domain

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Deposited on : 2018-07-19

Resolution : 2.90 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 (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 : 1.8.4, CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

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

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

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

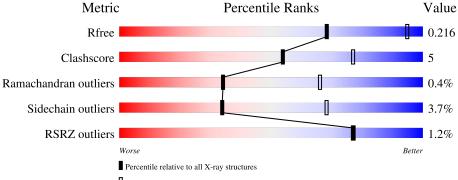
Validation Pipeline (wwPDB-VP) : 2.36

### 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 2.90 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.



Percentile relative to X-ray structures of similar resolution

Metric	Whole archive	Similar resolution
Metric	$(\#  ext{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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% 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	254	.%	80%	14% • 6%		
2	В	3	33%	67%			



# 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 1997 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 protein called Envelopment polyprotein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	A	240	Total 1958	C 1250	N 329	O 364	S 15	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	722	GLY	-	expression tag	UNP P04505
A	723	GLY	-	expression tag	UNP P04505
A	724	TRP	-	expression tag	UNP P04505
A	725	SER	_	expression tag	UNP P04505
A	726	HIS	-	expression tag	UNP P04505
A	727	PRO	_	expression tag	UNP P04505
A	728	GLN	-	expression tag	UNP P04505
Α	729	PHE	-	expression tag	UNP P04505
A	730	GLU	-	expression tag	UNP P04505
A	731	LYS	-	expression tag	UNP P04505

• Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[al pha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	В	3	Total 38	C 22	N 2	O 14	0	0	0

• Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).



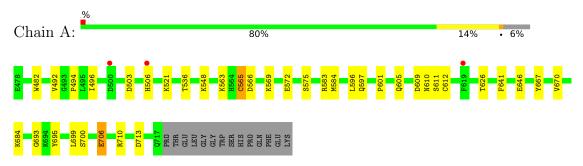
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0



# 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 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: Envelopment polyprotein



• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain B: 33% 67%





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants	125.44Å 125.44Å 46.55Å	Danagitan
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor
Resolution (Å)	46.55 - 2.90	Depositor
Resolution (A)	46.55 - 2.90	EDS
% Data completeness	99.7 (46.55-2.90)	Depositor
(in resolution range)	99.7 (46.55-2.90)	EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.58 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
D.D.	0.167 , 0.216	Depositor
$R, R_{free}$	0.167 , $0.216$	DCC
$R_{free}$ test set	945 reflections (9.91%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	95.4	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.34, 81.3	EDS
L-test for twinning <sup>2</sup>	$< L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.041 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	1997	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.42% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<sup>&</sup>lt;sup>1</sup>Intensities 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: FUC, NAG, CL

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).

Mal	Chain	Bond	lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.46	0/2003	0.62	2/2697 (0.1%)	

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$\operatorname{Ideal}(^{o})$
1	A	496	ILE	CB-CA-C	-5.06	101.47	111.60
1	A	565	CYS	CA-CB-SG	-5.02	104.97	114.00

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	1958	0	1919	20	0
2	В	38	0	34	2	0
3	A	1	0	0	0	0
All	All	1997	0	1953	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 5.

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



magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${\rm distance}({\rm \AA})$	overlap (Å)
1:A:612:CYS:HB2	1:A:684:LYS:HB3	1.75	0.69
2:B:2:NAG:H83	2:B:2:NAG:H3	1.85	0.57
1:A:646:GLU:OE2	1:A:710:ARG:NH2	2.39	0.56
1:A:503:ASP:HB3	1:A:506:HIS:ND1	2.22	0.55
1:A:548:LYS:NZ	1:A:713:ASP:O	2.38	0.55
1:A:492:VAL:HG23	1:A:494:PRO:HD2	1.95	0.49
1:A:569:LYS:O	1:A:572:GLU:HB2	2.12	0.49
1:A:482:TRP:CD1	1:A:521:LYS:HD2	2.47	0.48
1:A:597:GLN:OE1	1:A:626:THR:OG1	2.16	0.48
1:A:700:SER:OG	2:B:3:FUC:H3	2.13	0.48
1:A:667:TYR:CD1	1:A:670:VAL:HG21	2.50	0.45
1:A:641:PHE:CE1	1:A:693:GLY:HA3	2.54	0.43
1:A:706:GLU:OE1	1:A:706:GLU:N	2.43	0.43
1:A:695:TYR:CZ	1:A:699:LEU:HD21	2.55	0.42
1:A:601:PRO:O	1:A:605:GLN:HB2	2.20	0.42
1:A:605:GLN:O	1:A:609:ASP:HB2	2.19	0.42
1:A:667:TYR:CE1	1:A:670:VAL:HG21	2.56	0.41
1:A:506:HIS:ND1	1:A:506:HIS:N	2.69	0.41
1:A:566:ASP:HA	1:A:569:LYS:HB3	2.02	0.41
1:A:596:LEU:HD23	1:A:596:LEU:HA	1.66	0.40
1:A:548:LYS:HD3	1:A:548:LYS:HA	1.92	0.40

There are no symmetry-related clashes.

#### 5.3 Torsion angles (i)

#### 5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	238/254 (94%)	232 (98%)	5 (2%)	1 (0%)	34 66

#### All (1) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	611	SER

#### 5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	215/227~(95%)	207 (96%)	8 (4%)	34 68	

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	536	THR
1	A	563	LYS
1	A	565	CYS
1	A	575	SER
1	A	583	ARG
1	A	584	MET
1	A	610	ASN
1	A	706	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

#### 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)

3 monosaccharides 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 Chair		Chain	Chain Res Lii		Bo	ond leng	$ ag{ths}$	Bond angles		
MIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	В	1	2,1	14,14,15	0.79	1 (7%)	17,19,21	0.98	1 (5%)
2	NAG	В	2	2	14,14,15	0.77	1 (7%)	17,19,21	1.26	1 (5%)
2	FUC	В	3	2	10,10,11	1.24	1 (10%)	14,14,16	1.50	3 (21%)

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
2	NAG	В	1	2,1	-	3/6/23/26	0/1/1/1
2	NAG	В	2	2	-	4/6/23/26	0/1/1/1
2	FUC	В	3	2	-	-	0/1/1/1

#### All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\mathring{\mathbf{A}})$	$\operatorname{Ideal}(\text{\AA})$
2	В	1	NAG	O5-C1	2.67	1.48	1.43
2	В	3	FUC	O5-C5	2.57	1.49	1.43
2	В	2	NAG	O5-C1	2.05	1.47	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	В	2	NAG	C2-N2-C7	4.12	128.76	122.90
2	В	3	FUC	C1-C2-C3	3.53	114.00	109.67
2	В	1	NAG	C1-O5-C5	3.14	116.45	112.19
2	В	3	FUC	O5-C5-C4	2.59	114.17	109.52
2	В	3	FUC	C1-O5-C5	2.19	117.73	112.78

There are no chirality outliers.

All (7) torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
2	В	1	NAG	O5-C5-C6-O6
2	В	2	NAG	C8-C7-N2-C2
2	В	2	NAG	O7-C7-N2-C2
2	В	1	NAG	C4-C5-C6-O6
2	В	2	NAG	O5-C5-C6-O6
2	В	1	NAG	C3-C2-N2-C7
2	В	2	NAG	C3-C2-N2-C7

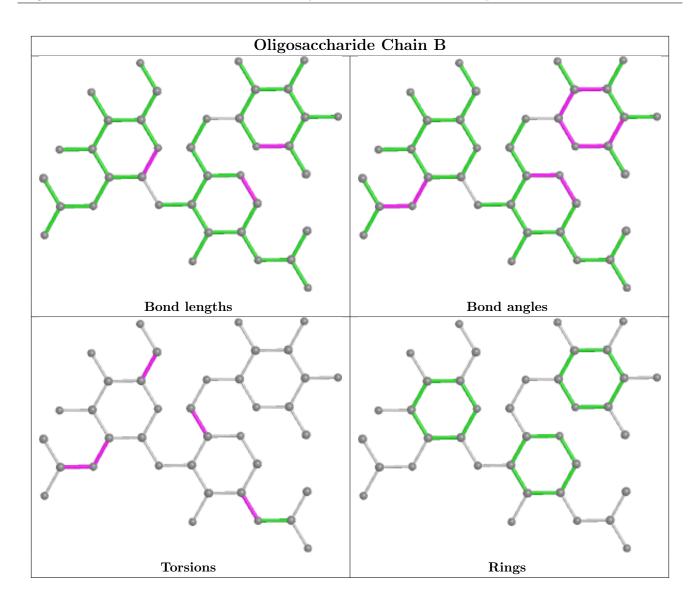
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	3	FUC	1	0
2	В	2	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





### 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 (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>	$\# \mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q < 0.9
1	A	240/254 (94%)	-0.09	3 (1%) 77 77	66, 96, 155, 222	0

All (3) RSRZ outliers are listed below:

Mol	Chain	$\operatorname{Res}$	Type	RSRZ
1	A	500	ASP	3.9
1	A	619	PHE	2.2
1	A	506	HIS	2.2

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

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

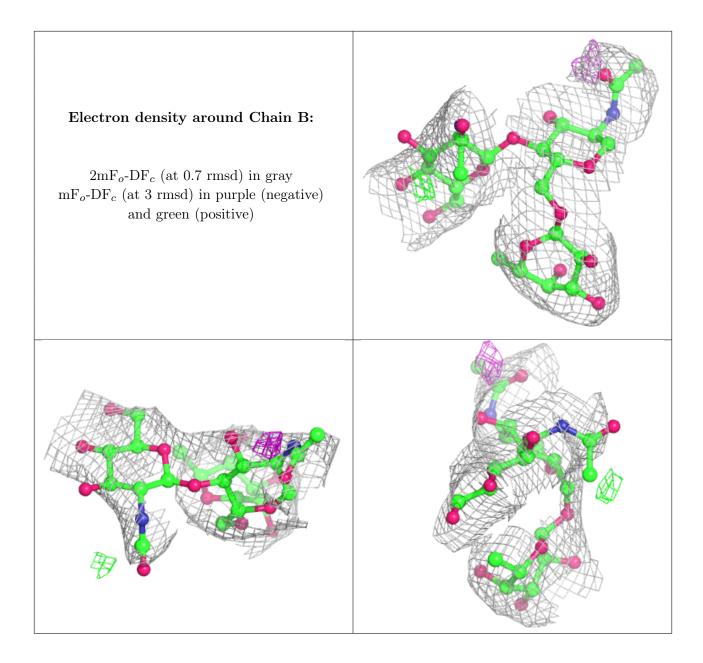
### 6.3 Carbohydrates (i)

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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
2	NAG	В	2	14/15	0.86	0.20	133,141,158,160	0
2	FUC	В	3	10/11	0.93	0.16	113,129,143,152	0
2	NAG	В	1	14/15	0.94	0.13	119,134,147,151	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.





### 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	$\operatorname{Res}$	Atoms	RSCC	RSR	$B$ -factors $(A^2)$	Q<0.9
3	CL	A	804	1/1	0.94	0.19	94,94,94,94	0



# 6.5 Other polymers (i)

There are no such residues in this entry.

