

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

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PDB ID : 4J4W

Title : Crystal structure of BueVN

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Deposited on : 2013-02-07

Resolution : 2.66 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 following versions of software and data (see references (1)) were used in the production of this report:

 $\begin{array}{ccc} Mol Probity & : & 4.02b\text{-}467 \\ Xtriage & (Phenix) & : & 1.13 \end{array}$ 

EDS : 2.11

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

Refmac: 5.8.0158

 $\begin{array}{cccc} & CCP4 & : & 7.0.044 \; (Gargrove) \\ Ideal \; geometry \; (proteins) & : & Engh \; \& \; Huber \; (2001) \end{array}$ 

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

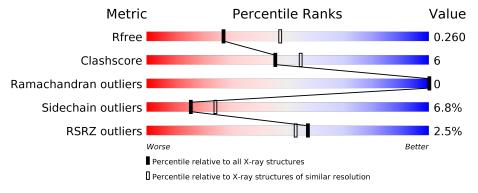
Validation Pipeline (wwPDB-VP) : 2.11

# 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 2.66 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
$R_{free}$	130704	$1332 \ (2.68-2.64)$
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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			
			2%			
1	Α	246	80%	17%	••	



# 2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1901 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 Nucleocapsid.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	243	Total	С	N	О	S	0	0	0
1	A	Z40	1871	1184	325	351	11	0	U	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference	
A	-2	SER	_	EXPRESSION TAG	UNP A7KCP4	
A	-1	ASN	-	EXPRESSION TAG	UNP A7KCP4	
A	0	ALA	_	EXPRESSION TAG	UNP A7KCP4	

• Molecule 2 is water.

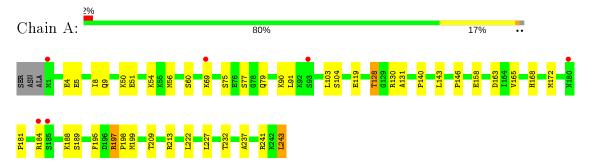
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	30	Total O 30 30	0	0



# 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: Nucleocapsid





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants	97.11Å 97.11Å 176.64Å	Danagitan
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.67 - 2.66	Depositor
Resolution (A)	32.67 - 2.66	EDS
% Data completeness	100.0 (32.67-2.66)	Depositor
(in resolution range)	100.0 (32.67-2.66)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	9.23 (at 2.65Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
D D	0.211 , 0.254	Depositor
$R, R_{free}$	0.216 , 0.260	DCC
$R_{free}$ test set	604 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.9	Xtriage
Anisotropy	0.148	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.37, 38.3	EDS
L-test for twinning <sup>2</sup>	$< L >=0.49, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	1901	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.00% 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)

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		
MIOI		RMSZ	# Z >5	RMSZ	# Z  > 5	
1	A	0.47	0/1906	0.55	0/2577	

There are no bond length outliers.

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	1871	0	1904	24	1
2	A	30	0	0	1	0
All	All	1901	0	1904	24	1

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 6.

All (24) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{l}  ext{Interatomic} \  ext{distance} \ ( ext{Å}) \end{array}$	Clash overlap (Å)	
1:A:90:LYS:O	2:A:318:HOH:O	1.90	0.87	
1:A:165:VAL:HG13	1:A:199:MET:HE2	1.60	0.84	
1:A:128:THR:HG22	1:A:131:ALA:H	1.40	0.83	
1:A:56:MET:HE2	1:A:91:LEU:HD11	1.82	0.62	

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A 4 1	A 4 0	Interatomic	Clash	
Atom-1	Atom-2	${f distance} \; ({f \AA})$	$overlap( ext{Å})$	
1:A:128:THR:CG2	1:A:131:ALA:H	2.13	0.56	
1:A:75:SER:O	1:A:79:GLN:HG3	2.05	0.56	
1:A:181:PRO:HA	1:A:184:ARG:HG3	1.89	0.55	
1:A:197:ARG:HG3	1:A:198:PRO:HD3	1.94	0.49	
1:A:168:HIS:HD2	1:A:199:MET:CE	2.26	0.49	
1:A:60:SER:HB2	1:A:103:LEU:HB3	1.96	0.48	
1:A:4:GLU:O	1:A:8:ILE:HG12	2.12	0.48	
1:A:128:THR:HG22	1:A:131:ALA:N	2.20	0.47	
1:A:209:THR:O	1:A:213:ARG:HG3	2.15	0.47	
1:A:128:THR:HG23	1:A:130:ARG:H	1.80	0.47	
1:A:237:ALA:O	1:A:241:ARG:HG3	2.15	0.47	
1:A:146:PRO:HG3	1:A:172:MET:HE1	1.97	0.46	
1:A:5:GLU:O	1:A:9:GLN:HG2	2.15	0.46	
1:A:195:PHE:O	1:A:198:PRO:HD2	2.16	0.45	
1:A:168:HIS:HD2	1:A:199:MET:HE3	1.82	0.44	
1:A:128:THR:HG23	1:A:130:ARG:N	2.33	0.43	
1:A:50:LYS:O	1:A:54:LYS:HG3	2.19	0.42	
1:A:140:PRO:HG2	1:A:143:LEU:HD12	2.01	0.42	
1:A:91:LEU:HA	1:A:91:LEU:HD23	1.81	0.41	
1:A:163:ASP:OD1	1:A:241:ARG:NH2	2.49	0.40	

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$egin{aligned}  ext{Interatomic} \  ext{distance} \ ( ext{Å}) \end{aligned}$	$egin{array}{c} { m Clash} \\ { m overlap} \ ({ m \AA}) \end{array}$
1:A:243:LEU:CG	1:A:243:LEU:CD1[5 655]	1.54	0.66

### 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	241/246 (98%)	230 (95%)	11 (5%)	0	100 100



There are no Ramachandran outliers to report.

#### 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	$206/208 \; (99\%)$	192 (93%)	14 (7%)	16 24

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	GLU
1	A	69	LYS
1	A	77	SER
1	A	104	SER
1	A	119	GLU
1	A	128	THR
1	A	158	GLU
1	A	188	LYS
1	A	189	SER
1	A	197	ARG
1	A	222	LEU
1	A	227	LEU
1	A	232	THR
1	A	243	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	HIS

#### 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,  $95^{th}$  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	$\langle { m RSRZ} \rangle$	#RS	$\mathbf{SRZ}$	>2	$OWAB(\AA^2)$	Q < 0.9
1	A	243/246 (98%)	-0.03	6 (2%)	57	53	36, 54, 73, 81	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	3.7
1	A	185	SER	3.7
1	A	184	ARG	2.9
1	A	69	LYS	2.8
1	A	180	ASN	2.5
1	A	93	SER	2.3

## 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.

