



# wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 25, 2026 – 12:36 PM EDT

PDB ID : 2YIB / pdb\_00002yib  
Title : Structure of the RNA polymerase VP1 from Infectious Pancreatic Necrosis Virus  
Authors : Graham, S.C.; Sarin, L.P.; Bahar, M.W.; Myers, R.A.; Stuart, D.I.; Bamford, D.H.; Grimes, J.M.  
Deposited on : 2011-05-11  
Resolution : 3.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

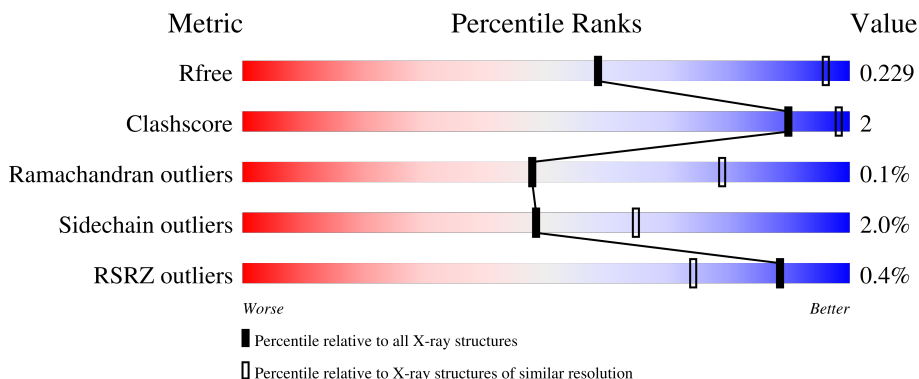
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.80 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))
$R_{free}$	180053	1065 (3.96-3.64)
Clashscore	190562	1012 (3.94-3.66)
Ramachandran outliers	187476	1048 (3.96-3.64)
Sidechain outliers	187428	1043 (3.96-3.64)
RSRZ outliers	180081	1064 (3.96-3.64)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	853	 86% 6% • 8%
1	B	853	 85% 7% 8%
1	C	853	 84% 8% 8%
2	D	770	 % 92% 7% •

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 24183 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 RNA-DIRECTED RNA POLYMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	789	6159	3908	1038	1187	26	0	0	0
1	B	789	6155	3906	1038	1185	26	0	0	0
1	C	789	6159	3908	1038	1187	26	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	846	VAL	-	expression tag	UNP P22173
A	847	GLU	-	expression tag	UNP P22173
A	848	HIS	-	expression tag	UNP P22173
A	849	HIS	-	expression tag	UNP P22173
A	850	HIS	-	expression tag	UNP P22173
A	851	HIS	-	expression tag	UNP P22173
A	852	HIS	-	expression tag	UNP P22173
A	853	HIS	-	expression tag	UNP P22173
B	846	VAL	-	expression tag	UNP P22173
B	847	GLU	-	expression tag	UNP P22173
B	848	HIS	-	expression tag	UNP P22173
B	849	HIS	-	expression tag	UNP P22173
B	850	HIS	-	expression tag	UNP P22173
B	851	HIS	-	expression tag	UNP P22173
B	852	HIS	-	expression tag	UNP P22173
B	853	HIS	-	expression tag	UNP P22173
C	846	VAL	-	expression tag	UNP P22173
C	847	GLU	-	expression tag	UNP P22173
C	848	HIS	-	expression tag	UNP P22173
C	849	HIS	-	expression tag	UNP P22173
C	850	HIS	-	expression tag	UNP P22173
C	851	HIS	-	expression tag	UNP P22173
C	852	HIS	-	expression tag	UNP P22173

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Chain	Residue	Modelled	Actual	Comment	Reference
C	853	HIS	-	expression tag	UNP P22173

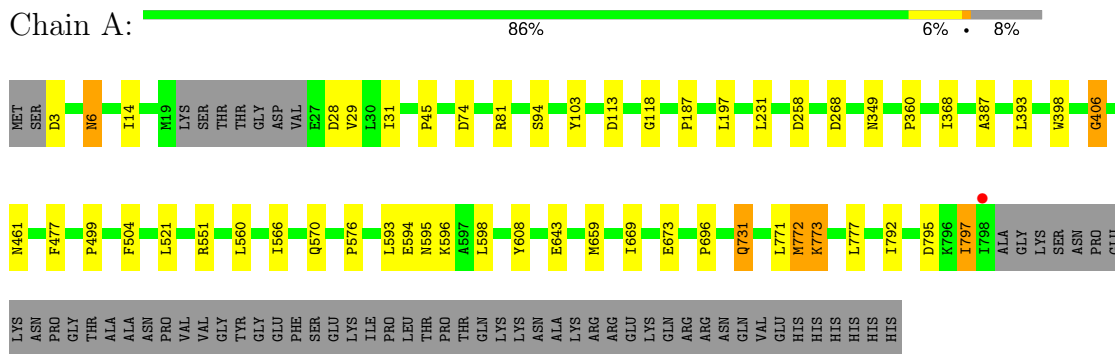
- Molecule 2 is a protein called RNA-DIRECTED RNA POLYMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	D	761	5710	3622	966	1097	25	415	0	0

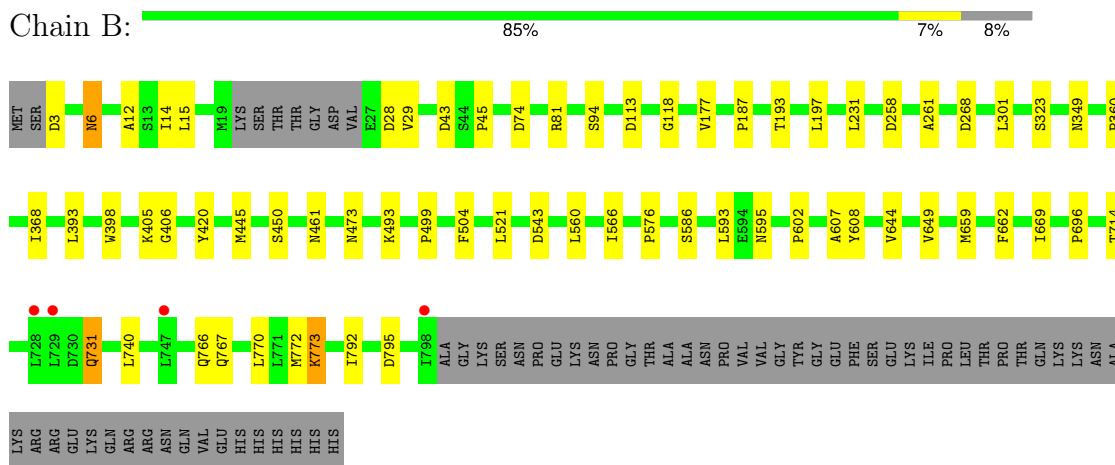
### 3 Residue-property plots

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.

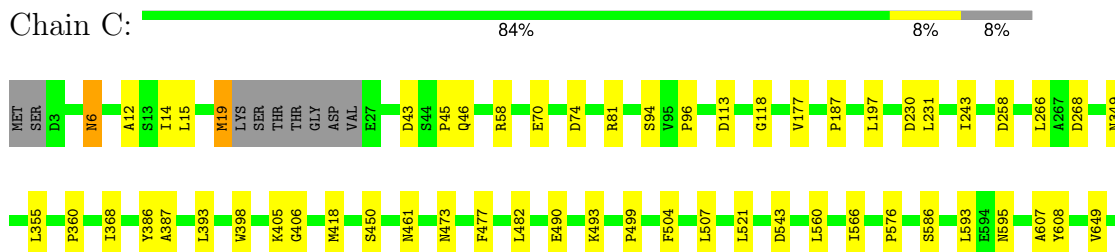
- Molecule 1: RNA-DIRECTED RNA POLYMERASE

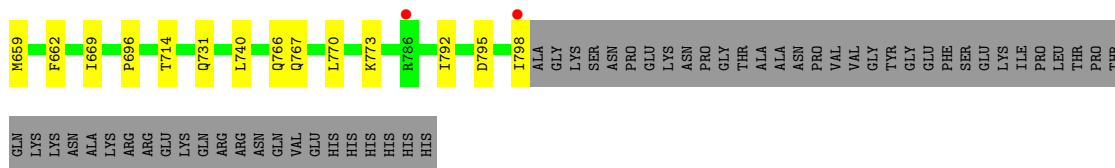


- Molecule 1: RNA-DIRECTED RNA POLYMERASE

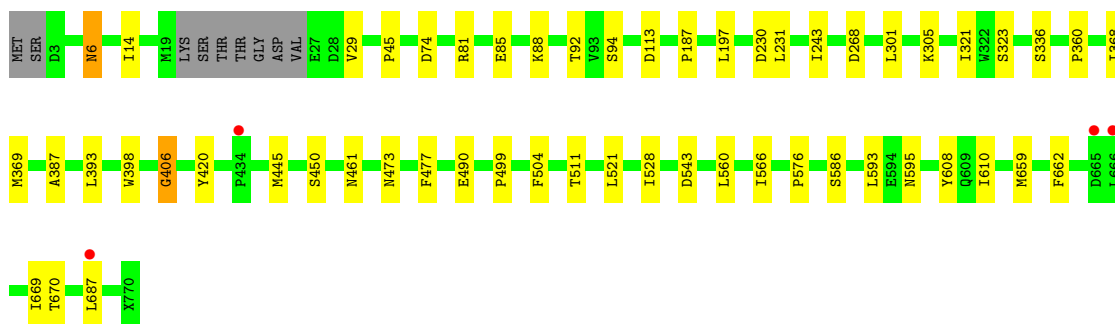


- Molecule 1: RNA-DIRECTED RNA POLYMERASE





● Molecule 2: RNA-DIRECTED RNA POLYMERASE



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	119.74Å 195.45Å 197.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.43 – 3.80 44.43 – 3.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (44.43-3.80) 98.0 (44.43-3.80)	Depositor EDS
$R_{merge}$	0.30	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.91 (at 3.77Å)	Xtrriage
Refinement program	BUSTER 2.9.2	Depositor
R, $R_{free}$	0.184 , 0.213 0.203 , 0.229	Depositor DCC
$R_{free}$ test set	2307 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	88.6	Xtrriage
Anisotropy	0.182	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 115.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.41$ , $\langle L^2 \rangle = 0.23$	Xtrriage
Estimated twinning fraction	0.056 for -h,l,k	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	24183	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	99.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.98% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.79	0/6296	1.25	13/8573 (0.2%)
1	B	0.77	0/6292	1.31	22/8568 (0.3%)
1	C	0.81	0/6296	1.31	18/8573 (0.2%)
2	D	0.79	0/5420	1.29	14/7383 (0.2%)
All	All	0.79	0/24304	1.29	67/33097 (0.2%)

There are no bond length outliers.

The worst 5 of 67 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	773	LYS	N-CA-C	-7.43	104.19	113.55
1	C	773	LYS	N-CA-C	-7.26	104.28	113.43
1	B	118	GLY	CA-C-N	6.90	129.84	120.38
1	B	118	GLY	C-N-CA	6.90	129.84	120.38
1	A	773	LYS	N-CA-C	-6.41	105.38	113.72

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	6159	0	6090	24	0
1	B	6155	0	6086	26	0
1	C	6159	0	6090	28	0
2	D	5710	0	5296	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	24183	0	23562	87	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 2.

The worst 5 of 87 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:644:VAL:HG22	2:D:92:THR:CG2	1.90	1.01
1:A:731:GLN:HG2	1:A:772:MET:HE1	1.50	0.93
1:B:644:VAL:HG22	2:D:92:THR:HG21	1.54	0.88
1:B:644:VAL:HG22	2:D:92:THR:HG23	1.61	0.81
1:C:187:PRO:HB3	1:C:197:LEU:HD11	1.73	0.70

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	785/853 (92%)	761 (97%)	23 (3%)	1 (0%)	48 79
1	B	785/853 (92%)	761 (97%)	23 (3%)	1 (0%)	48 79
1	C	785/853 (92%)	758 (97%)	26 (3%)	1 (0%)	48 79
2	D	674/770 (88%)	651 (97%)	22 (3%)	1 (0%)	48 79
All	All	3029/3329 (91%)	2931 (97%)	94 (3%)	4 (0%)	48 79

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	406	GLY

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Mol	Chain	Res	Type
1	B	406	GLY
1	C	406	GLY
1	A	406	GLY

### 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	669/739 (90%)	656 (98%)	13 (2%)	50	66
1	B	668/739 (90%)	654 (98%)	14 (2%)	47	64
1	C	669/739 (90%)	655 (98%)	14 (2%)	47	64
2	D	573/593 (97%)	562 (98%)	11 (2%)	50	66
All	All	2579/2810 (92%)	2527 (98%)	52 (2%)	48	65

5 of 52 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	14	ILE
1	C	450	SER
2	D	586	SER
1	C	15	LEU
1	C	94	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	429	ASN
1	C	793	GLN
1	C	768	GLN
2	D	226	ASN
1	B	429	ASN

### 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 oligosaccharides 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	D	3

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	687:LEU	C	688:UNK	N	3.36
1	D	767:UNK	C	768:UNK	N	1.03
1	D	700:UNK	C	701:UNK	N	1.02

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	789/853 (92%)	-0.42	1 (0%) 92 87	50, 80, 118, 174	0
1	B	789/853 (92%)	-0.22	4 (0%) 87 71	62, 113, 183, 214	0
1	C	789/853 (92%)	-0.45	2 (0%) 90 78	49, 85, 131, 179	0
2	D	678/770 (88%)	-0.25	4 (0%) 85 68	58, 103, 162, 190	0
All	All	3045/3329 (91%)	-0.34	11 (0%) 88 74	49, 93, 158, 214	0

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	798	ILE	4.2
2	D	434	PRO	3.5
1	B	728	LEU	3.4
1	B	747	LEU	3.2
2	D	666	LEU	2.5

### 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 oligosaccharides 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.