



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 28, 2017 – 03:55 AM EDT

PDB ID : 5UKB  
Title : VSV N PROTEIN IN COMPLEX WITH INHIBITORY NANOBODY 1004  
Authors : Hanke, L.; Knockenhauer, K.E.; Ploegh, H.L.; Schwartz, T.U.  
Deposited on : unknown  
Resolution : 5.47 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

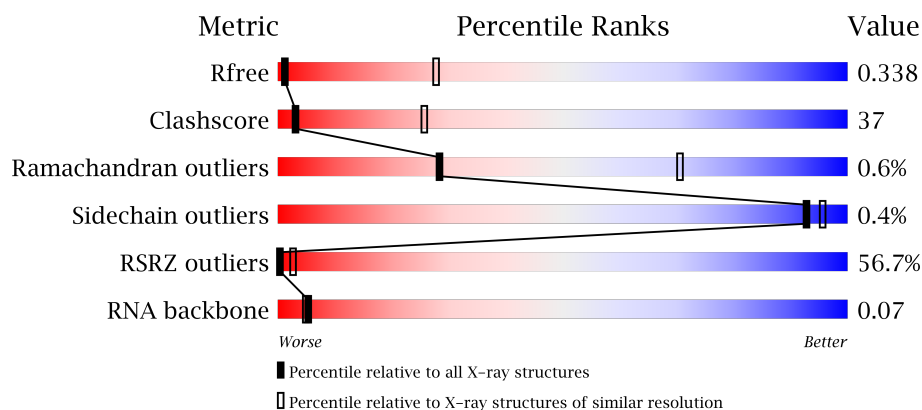
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 5.47 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))
$R_{free}$	100719	1052 (7.20-3.70)
Clashscore	112137	1021 (7.20-3.76)
Ramachandran outliers	110173	1082 (7.20-3.70)
Sidechain outliers	110143	1055 (7.20-3.70)
RSRZ outliers	101464	1061 (7.20-3.70)
RNA backbone	2435	1049 (7.80-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	138	<div> <div>50%</div> <div>84%</div> <div>12%</div> </div>
1	b	138	<div> <div>47%</div> <div>86%</div> <div>12%</div> </div>
1	c	138	<div> <div>65%</div> <div>84%</div> <div>12%</div> </div>
1	d	138	<div> <div>54%</div> <div>86%</div> <div>12%</div> </div>

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Mol	Chain	Length	Quality of chain
1	e	138	<div><div></div><div>66%</div><div>86%</div><div>12%</div></div>
2	A	423	<div><div></div><div>48%</div><div>73%</div><div>25%</div></div>
2	B	423	<div><div></div><div>56%</div><div>73%</div><div>26%</div></div>
2	C	423	<div><div></div><div>47%</div><div>74%</div><div>24%</div></div>
2	D	423	<div><div></div><div>61%</div><div>74%</div><div>24%</div></div>
2	E	423	<div><div></div><div>53%</div><div>74%</div><div>25%</div></div>
3	R	45	<div><div></div><div>100%</div><div>22%</div><div>56%</div><div>18%</div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 22210 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 Anti-vesicular stomatitis virus N VHH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	d	122	Total	C	N	O	S	0	0	0
			946	596	166	181	3			
1	c	122	Total	C	N	O	S	0	0	0
			946	596	166	181	3			
1	b	122	Total	C	N	O	S	0	0	0
			946	596	166	181	3			
1	a	122	Total	C	N	O	S	0	0	0
			946	596	166	181	3			
1	e	122	Total	C	N	O	S	0	0	0
			946	596	166	181	3			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
d	1	GLN	-	expression tag	UNP A0A192B6J5
d	2	VAL	-	expression tag	UNP A0A192B6J5
d	125	GLY	-	expression tag	UNP A0A192B6J5
d	126	GLY	-	expression tag	UNP A0A192B6J5
d	127	LEU	-	expression tag	UNP A0A192B6J5
d	128	PRO	-	expression tag	UNP A0A192B6J5
d	129	GLU	-	expression tag	UNP A0A192B6J5
d	130	THR	-	expression tag	UNP A0A192B6J5
d	131	GLY	-	expression tag	UNP A0A192B6J5
d	132	GLY	-	expression tag	UNP A0A192B6J5
d	133	HIS	-	expression tag	UNP A0A192B6J5
d	134	HIS	-	expression tag	UNP A0A192B6J5
d	135	HIS	-	expression tag	UNP A0A192B6J5
d	136	HIS	-	expression tag	UNP A0A192B6J5
d	137	HIS	-	expression tag	UNP A0A192B6J5
d	138	HIS	-	expression tag	UNP A0A192B6J5
c	1	GLN	-	expression tag	UNP A0A192B6J5
c	2	VAL	-	expression tag	UNP A0A192B6J5
c	125	GLY	-	expression tag	UNP A0A192B6J5

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Chain	Residue	Modelled	Actual	Comment	Reference
c	126	GLY	-	expression tag	UNP A0A192B6J5
c	127	LEU	-	expression tag	UNP A0A192B6J5
c	128	PRO	-	expression tag	UNP A0A192B6J5
c	129	GLU	-	expression tag	UNP A0A192B6J5
c	130	THR	-	expression tag	UNP A0A192B6J5
c	131	GLY	-	expression tag	UNP A0A192B6J5
c	132	GLY	-	expression tag	UNP A0A192B6J5
c	133	HIS	-	expression tag	UNP A0A192B6J5
c	134	HIS	-	expression tag	UNP A0A192B6J5
c	135	HIS	-	expression tag	UNP A0A192B6J5
c	136	HIS	-	expression tag	UNP A0A192B6J5
c	137	HIS	-	expression tag	UNP A0A192B6J5
c	138	HIS	-	expression tag	UNP A0A192B6J5
b	1	GLN	-	expression tag	UNP A0A192B6J5
b	2	VAL	-	expression tag	UNP A0A192B6J5
b	125	GLY	-	expression tag	UNP A0A192B6J5
b	126	GLY	-	expression tag	UNP A0A192B6J5
b	127	LEU	-	expression tag	UNP A0A192B6J5
b	128	PRO	-	expression tag	UNP A0A192B6J5
b	129	GLU	-	expression tag	UNP A0A192B6J5
b	130	THR	-	expression tag	UNP A0A192B6J5
b	131	GLY	-	expression tag	UNP A0A192B6J5
b	132	GLY	-	expression tag	UNP A0A192B6J5
b	133	HIS	-	expression tag	UNP A0A192B6J5
b	134	HIS	-	expression tag	UNP A0A192B6J5
b	135	HIS	-	expression tag	UNP A0A192B6J5
b	136	HIS	-	expression tag	UNP A0A192B6J5
b	137	HIS	-	expression tag	UNP A0A192B6J5
b	138	HIS	-	expression tag	UNP A0A192B6J5
a	1	GLN	-	expression tag	UNP A0A192B6J5
a	2	VAL	-	expression tag	UNP A0A192B6J5
a	125	GLY	-	expression tag	UNP A0A192B6J5
a	126	GLY	-	expression tag	UNP A0A192B6J5
a	127	LEU	-	expression tag	UNP A0A192B6J5
a	128	PRO	-	expression tag	UNP A0A192B6J5
a	129	GLU	-	expression tag	UNP A0A192B6J5
a	130	THR	-	expression tag	UNP A0A192B6J5
a	131	GLY	-	expression tag	UNP A0A192B6J5
a	132	GLY	-	expression tag	UNP A0A192B6J5
a	133	HIS	-	expression tag	UNP A0A192B6J5
a	134	HIS	-	expression tag	UNP A0A192B6J5
a	135	HIS	-	expression tag	UNP A0A192B6J5

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Chain	Residue	Modelled	Actual	Comment	Reference
a	136	HIS	-	expression tag	UNP A0A192B6J5
a	137	HIS	-	expression tag	UNP A0A192B6J5
a	138	HIS	-	expression tag	UNP A0A192B6J5
e	1	GLN	-	expression tag	UNP A0A192B6J5
e	2	VAL	-	expression tag	UNP A0A192B6J5
e	125	GLY	-	expression tag	UNP A0A192B6J5
e	126	GLY	-	expression tag	UNP A0A192B6J5
e	127	LEU	-	expression tag	UNP A0A192B6J5
e	128	PRO	-	expression tag	UNP A0A192B6J5
e	129	GLU	-	expression tag	UNP A0A192B6J5
e	130	THR	-	expression tag	UNP A0A192B6J5
e	131	GLY	-	expression tag	UNP A0A192B6J5
e	132	GLY	-	expression tag	UNP A0A192B6J5
e	133	HIS	-	expression tag	UNP A0A192B6J5
e	134	HIS	-	expression tag	UNP A0A192B6J5
e	135	HIS	-	expression tag	UNP A0A192B6J5
e	136	HIS	-	expression tag	UNP A0A192B6J5
e	137	HIS	-	expression tag	UNP A0A192B6J5
e	138	HIS	-	expression tag	UNP A0A192B6J5

- Molecule 2 is a protein called Nucleocapsid.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	419	Total	C	N	O	S	0	0	0
			3316	2111	556	631	18			
2	C	419	Total	C	N	O	S	0	0	0
			3316	2111	556	631	18			
2	B	419	Total	C	N	O	S	0	0	0
			3316	2111	556	631	18			
2	A	419	Total	C	N	O	S	0	0	0
			3316	2111	556	631	18			
2	E	419	Total	C	N	O	S	0	0	0
			3316	2111	556	631	18			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	MET	-	expression tag	UNP A6H4P1
D	1	ALA	-	expression tag	UNP A6H4P1
C	0	MET	-	expression tag	UNP A6H4P1
C	1	ALA	-	expression tag	UNP A6H4P1
B	0	MET	-	expression tag	UNP A6H4P1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1	ALA	-	expression tag	UNP A6H4P1
A	0	MET	-	expression tag	UNP A6H4P1
A	1	ALA	-	expression tag	UNP A6H4P1
E	0	MET	-	expression tag	UNP A6H4P1
E	1	ALA	-	expression tag	UNP A6H4P1

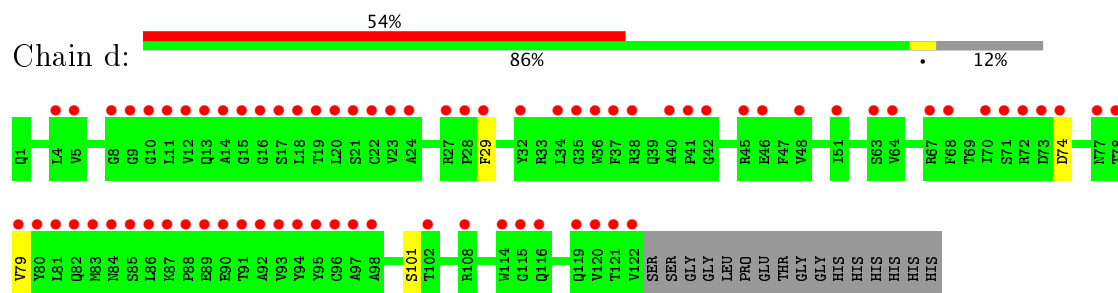
- Molecule 3 is a RNA chain called RNA (45-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	R	45	Total	C	N	O	P	0	0	0
			900	405	90	360	45			

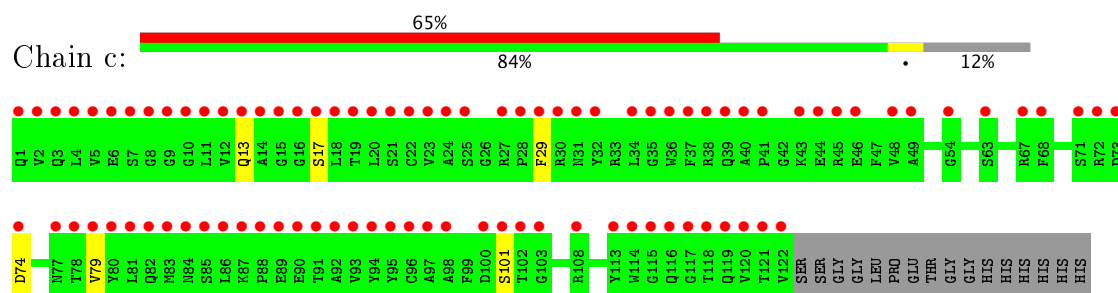
### 3 Residue-property plots

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.

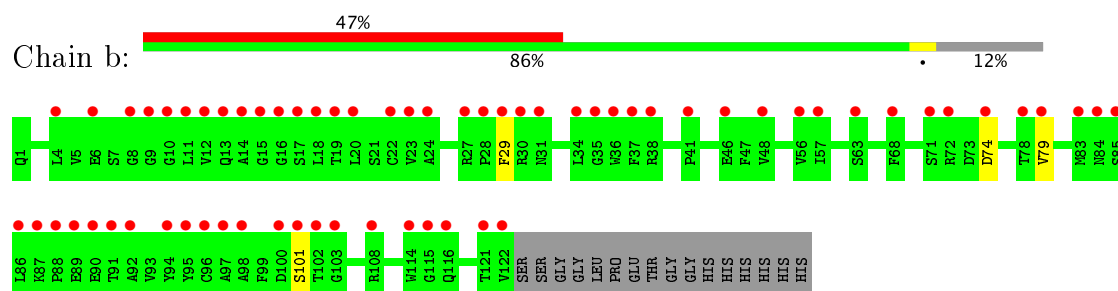
- Molecule 1: Anti-vesicular stomatitis virus N VHH



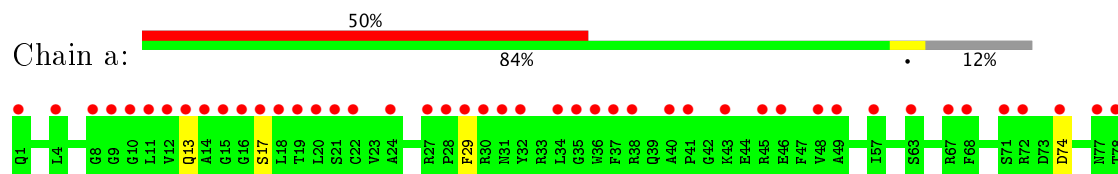
- Molecule 1: Anti-vesicular stomatitis virus N VHH



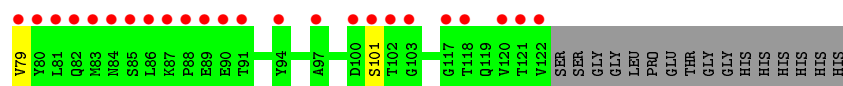
- Molecule 1: Anti-vesicular stomatitis virus N VHH



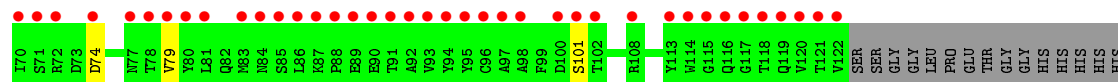
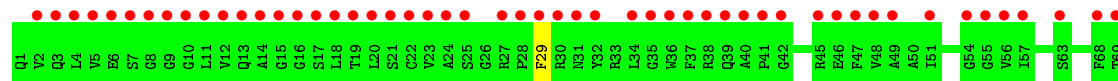
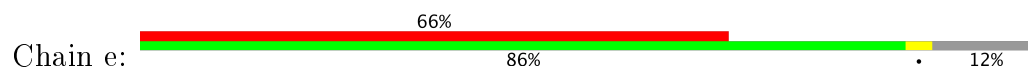
- Molecule 1: Anti-vesicular stomatitis virus N VHH





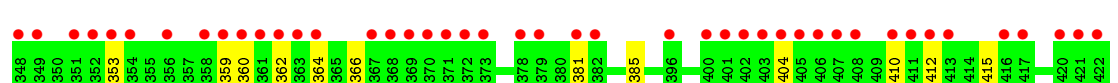
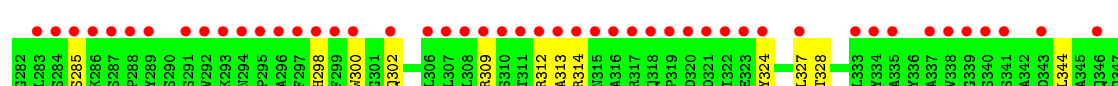
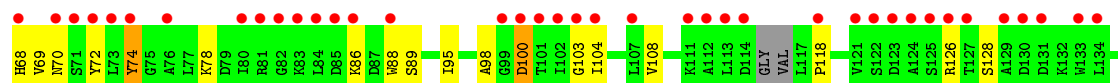
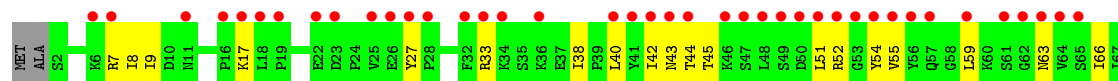
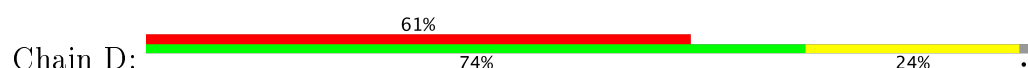


• Molecule 1: Anti-vesicular stomatitis virus N VHH

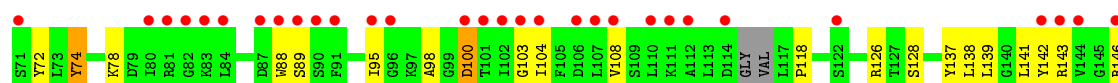


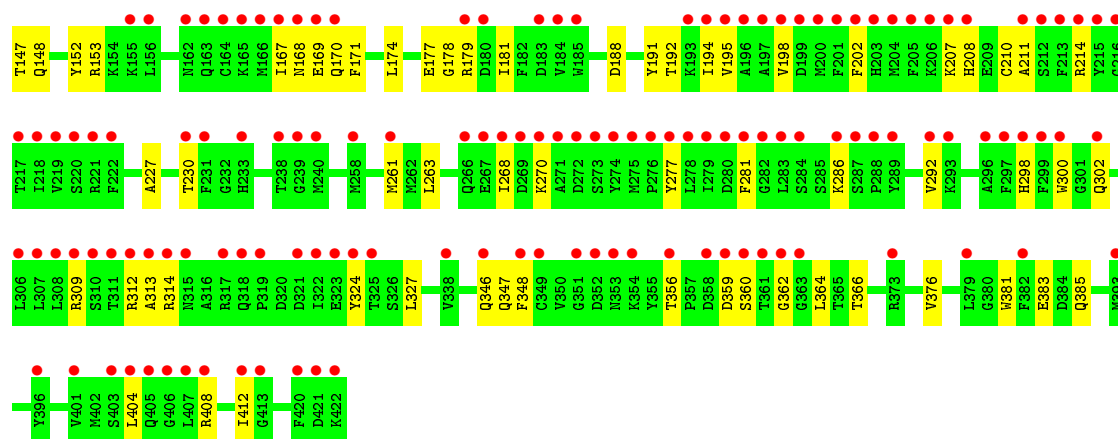
HIS

• Molecule 2: Nucleocapsid

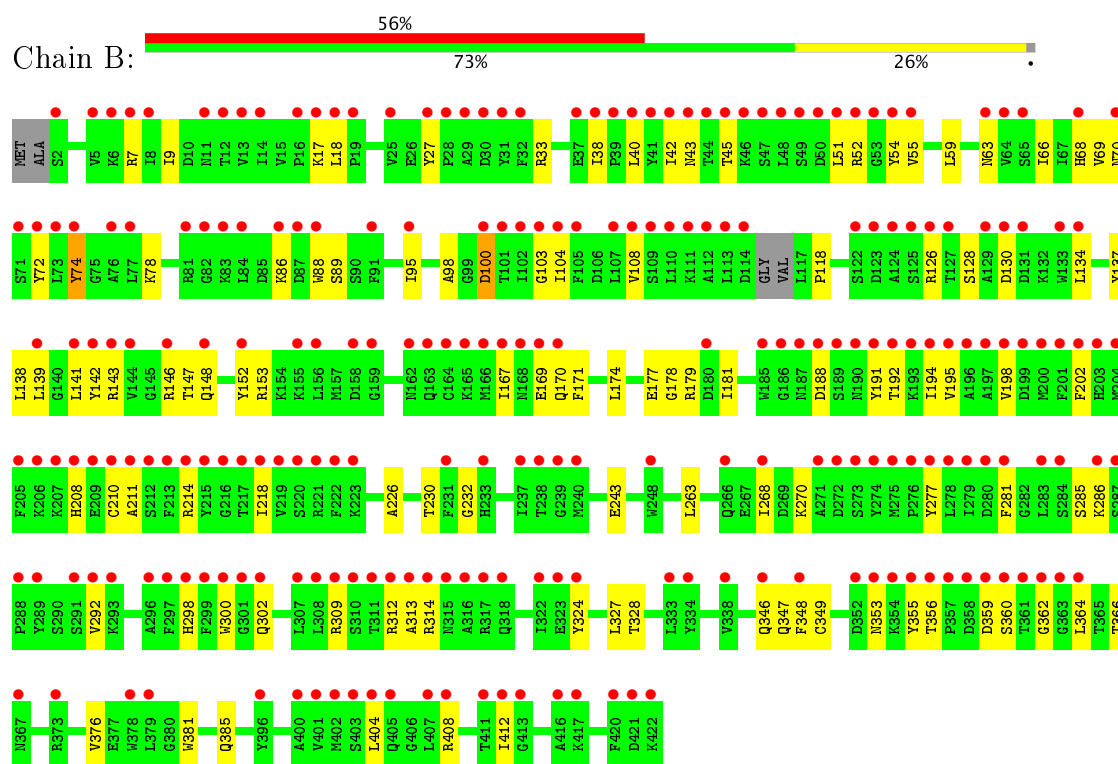


• Molecule 2: Nucleocapsid

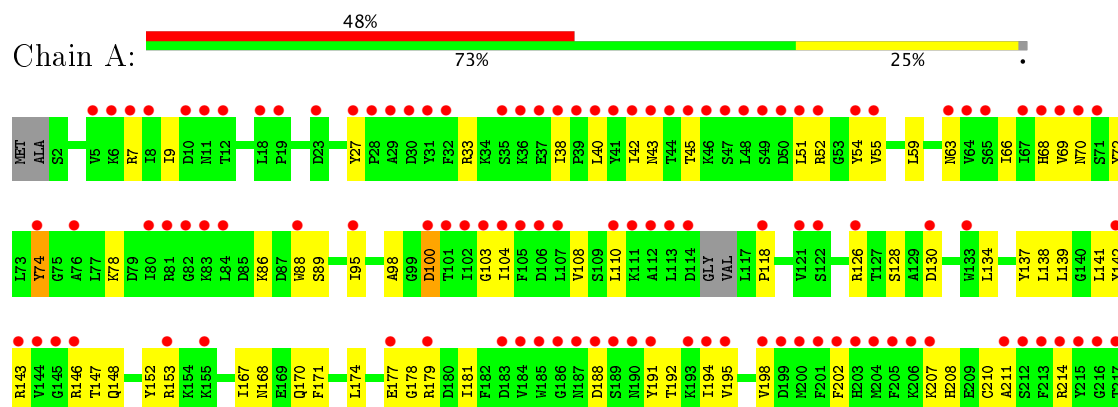


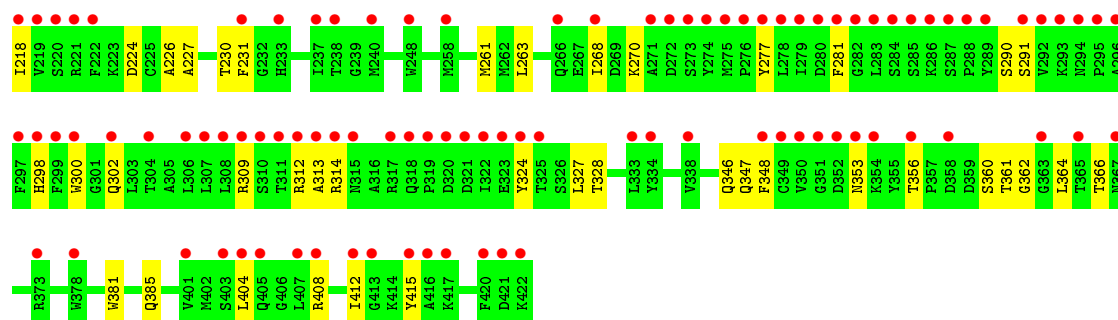


### • Molecule 2: Nucleocapsid

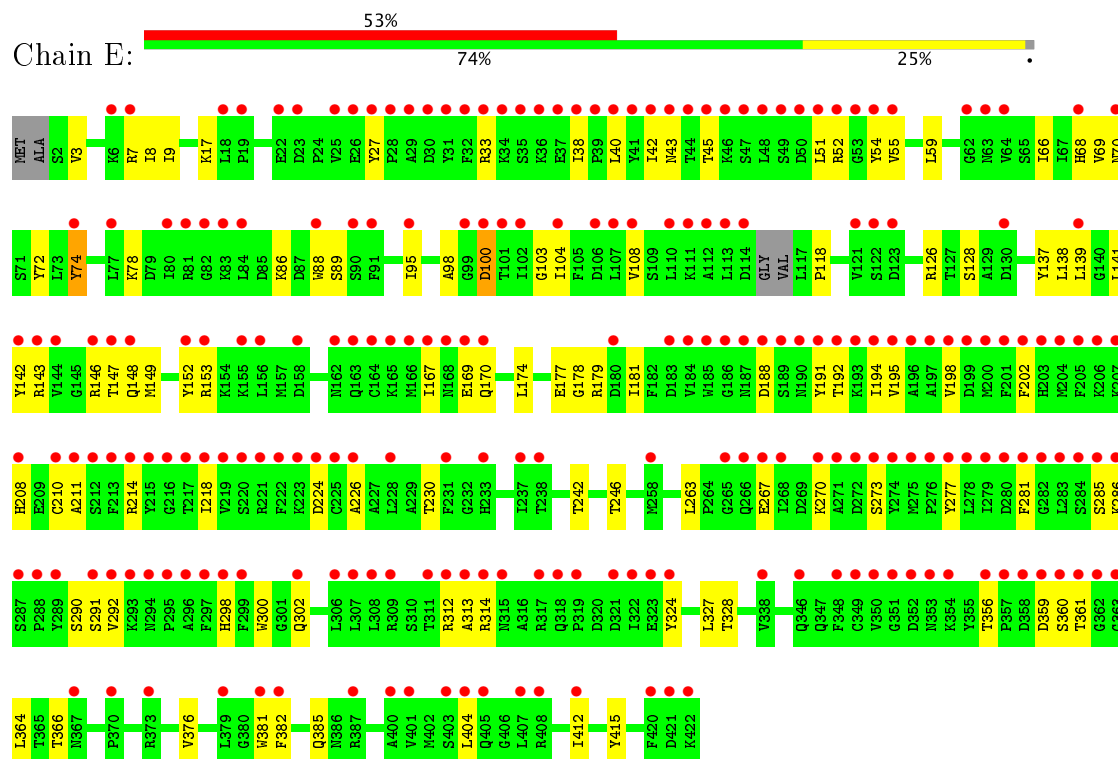


### • Molecule 2: Nucleocapsid

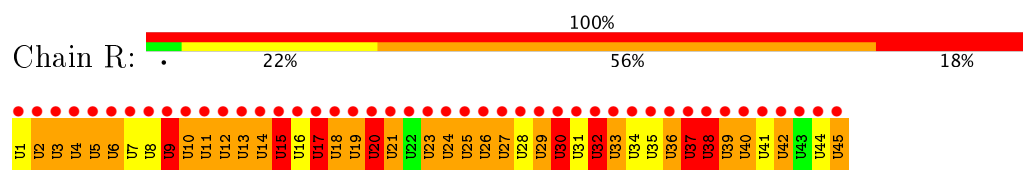




• Molecule 2: Nucleocapsid



• Molecule 3: RNA (45-MER)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	240.12Å 335.50Å 75.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	137.51 – 5.47 137.51 – 5.47	Depositor EDS
% Data completeness (in resolution range)	97.4 (137.51-5.47) 97.4 (137.51-5.47)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.17	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.24 (at 5.42Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, $R_{free}$	0.336 , 0.338 0.336 , 0.338	Depositor DCC
$R_{free}$ test set	1996 reflections (9.70%)	DCC
Wilson B-factor (Å <sup>2</sup> )	292.7	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 423.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.39$ , $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.81	EDS
Total number of atoms	22210	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	309.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.78% 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	0.25	0/967	0.55	1/1312 (0.1%)
1	b	0.25	0/967	0.55	1/1312 (0.1%)
1	c	0.26	0/967	0.55	1/1312 (0.1%)
1	d	0.25	0/967	0.55	1/1312 (0.1%)
1	e	0.26	0/967	0.55	1/1312 (0.1%)
2	A	0.24	0/3391	0.41	0/4589
2	B	0.24	0/3391	0.41	0/4589
2	C	0.24	0/3391	0.41	0/4589
2	D	0.24	0/3391	0.41	0/4589
2	E	0.24	0/3391	0.41	0/4589
3	R	0.94	0/989	1.56	10/1526 (0.7%)
All	All	0.31	0/22779	0.56	15/31031 (0.0%)

There are no bond length outliers.

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	R	20	U	P-O3'-C3'	8.95	130.44	119.70
3	R	9	U	O4'-C1'-N1	6.60	113.48	108.20
3	R	9	U	O4'-C4'-C3'	-6.11	97.89	104.00
3	R	34	U	O4'-C1'-N1	5.88	112.91	108.20
3	R	15	U	C2-N1-C1'	5.79	124.65	117.70

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	946	0	909	0	2
1	b	946	0	908	0	0
1	c	946	0	908	0	2
1	d	946	0	908	0	0
1	e	946	0	908	0	0
2	A	3316	0	3272	115	14
2	B	3316	0	3272	144	12
2	C	3316	0	3272	138	14
2	D	3316	0	3273	140	10
2	E	3316	0	3272	139	16
3	R	900	0	451	139	5
All	All	22210	0	21353	611	51

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 37.

The worst 5 of 611 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)
2:D:214:ARG:NH2	3:R:37:U:C5'	1.78	1.44
2:E:33:ARG:CD	2:E:103:GLY:O	26.07	1.43
2:C:33:ARG:CD	2:C:103:GLY:O	26.07	1.43
2:D:206:LYS:CE	3:R:38:U:OP1	1.65	1.42
2:C:33:ARG:CG	2:C:103:GLY:O	24.82	1.42

The worst 5 of 51 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	Interatomic distance (Å)	Clash overlap (Å)
2:C:169:GLU:CB	2:C:359:ASP:CB[1_554]	0.90	1.30
1:c:13:GLN:NE2	1:a:17:SER:O[4_477]	1.59	0.61
3:R:1:U:P	3:R:45:U:O3'[2_675]	1.60	0.60
2:D:360:SER:OG	2:E:177:GLU:OE1[1_556]	1.63	0.57
2:B:177:GLU:CD	2:A:360:SER:O[1_554]	1.67	0.53

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	120/138 (87%)	110 (92%)	8 (7%)	2 (2%)	11	51
1	b	120/138 (87%)	110 (92%)	8 (7%)	2 (2%)	11	51
1	c	120/138 (87%)	110 (92%)	8 (7%)	2 (2%)	11	51
1	d	120/138 (87%)	110 (92%)	8 (7%)	2 (2%)	11	51
1	e	120/138 (87%)	110 (92%)	8 (7%)	2 (2%)	11	51
2	A	415/423 (98%)	395 (95%)	19 (5%)	1 (0%)	51	84
2	B	415/423 (98%)	395 (95%)	19 (5%)	1 (0%)	51	84
2	C	415/423 (98%)	395 (95%)	19 (5%)	1 (0%)	51	84
2	D	415/423 (98%)	395 (95%)	19 (5%)	1 (0%)	51	84
2	E	415/423 (98%)	395 (95%)	19 (5%)	1 (0%)	51	84
All	All	2675/2805 (95%)	2525 (94%)	135 (5%)	15 (1%)	28	70

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	d	101	SER
1	c	101	SER
1	b	101	SER
1	a	101	SER
1	e	101	SER

### 5.3.2 Protein sidechains

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	98/110 (89%)	97 (99%)	1 (1%)	80	90
1	b	98/110 (89%)	97 (99%)	1 (1%)	80	90
1	c	98/110 (89%)	97 (99%)	1 (1%)	80	90
1	d	98/110 (89%)	97 (99%)	1 (1%)	80	90
1	e	98/110 (89%)	97 (99%)	1 (1%)	80	90
2	A	361/363 (99%)	360 (100%)	1 (0%)	94	96
2	B	361/363 (99%)	360 (100%)	1 (0%)	94	96
2	C	361/363 (99%)	360 (100%)	1 (0%)	94	96
2	D	361/363 (99%)	360 (100%)	1 (0%)	94	96
2	E	361/363 (99%)	360 (100%)	1 (0%)	94	96
All	All	2295/2365 (97%)	2285 (100%)	10 (0%)	93	95

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	b	29	PHE
2	B	100	ASP
2	A	100	ASP
2	C	100	ASP
1	a	29	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
1	b	77	ASN
1	b	82	GLN
2	A	208	HIS
2	C	208	HIS
1	a	82	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	R	44/45 (97%)	35 (79%)	8 (18%)

5 of 35 RNA backbone outliers are listed below:



Mol	Chain	Res	Type
3	R	2	U
3	R	3	U
3	R	4	U
3	R	5	U
3	R	6	U

5 of 8 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	R	12	U
3	R	39	U
3	R	20	U
3	R	6	U
3	R	14	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](#)

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 ⓘ

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	122/138 (88%)	3.73	69 (56%) 0 3	303, 317, 325, 340	0
1	b	122/138 (88%)	2.70	65 (53%) 0 3	346, 363, 371, 373	0
1	c	122/138 (88%)	5.12	90 (73%) 0 2	311, 328, 337, 348	0
1	d	122/138 (88%)	3.37	75 (61%) 0 2	371, 385, 392, 396	0
1	e	122/138 (88%)	4.75	91 (74%) 0 2	358, 378, 388, 390	0
2	A	419/423 (99%)	2.39	205 (48%) 0 3	236, 284, 336, 356	0
2	B	419/423 (99%)	2.86	235 (56%) 0 3	235, 288, 348, 365	0
2	C	419/423 (99%)	2.40	199 (47%) 0 3	242, 285, 342, 367	0
2	D	419/423 (99%)	3.07	258 (61%) 0 2	246, 309, 361, 377	0
2	E	419/423 (99%)	2.80	226 (53%) 0 3	247, 302, 355, 373	0
3	R	45/45 (100%)	7.78	45 (100%) 0 0	289, 313, 332, 338	0
All	All	2750/2850 (96%)	3.06	1558 (56%) 0 3	235, 311, 380, 396	0

The worst 5 of 1558 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	e	8	GLY	23.1
1	a	13	GLN	22.6
1	c	13	GLN	22.4
1	c	16	GLY	22.2
1	a	15	GLY	20.4

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

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.