wwPDB X-ray Structure Validation Summary Report ⓘ

May 23, 2020 – 12:39 am BST

PDB ID	:	4JNG
Title	:	Schmallenberg virus nucleoprotein-RNA complex
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Deposited on	:	2013-03-15
Resolution	:	2.12 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
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

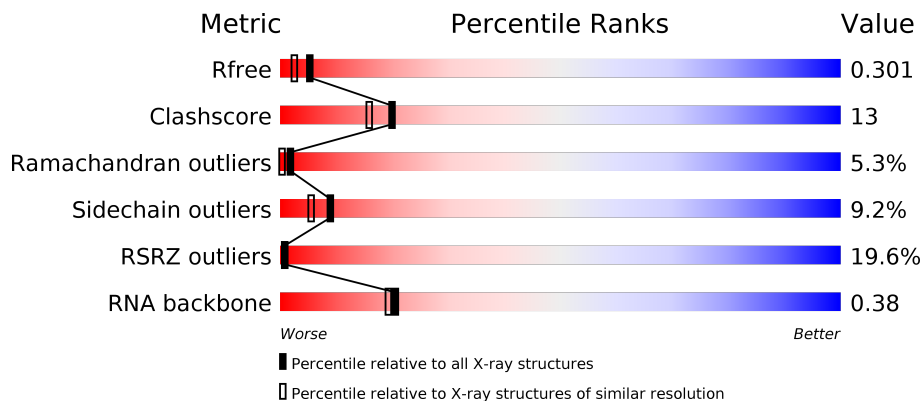|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.12 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.
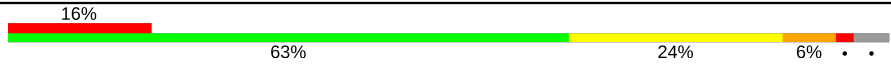
Metric	Percentile Ranks	Value
Rfree		0.301
Clashscore		13
Ramachandran outliers		5.3%
Sidechain outliers		9.2%
RSRZ outliers		19.6%
RNA backbone		0.38

Worse   *Better*

■ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6241 (2.14-2.10)
Clashscore	141614	6778 (2.14-2.10)
Ramachandran outliers	138981	6705 (2.14-2.10)
Sidechain outliers	138945	6706 (2.14-2.10)
RSRZ outliers	127900	6112 (2.14-2.10)
RNA backbone	3102	1013 (2.58-1.66)

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
1	L	42	52% 29% 19%
2	A	233	22% 76% 18% · ·
2	B	233	9% 69% 19% · · 8%
2	C	233	30% 64% 24% 7% · ·

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Mol	Chain	Length	Quality of chain
2	D	233	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8038 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 RNA (42-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	42	Total 840	C 378	N 84	O 336	P 42	0	0	0

- Molecule 2 is a protein called Nucleocapsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	226	Total 1796	C 1168	N 301	O 315	S 12	0	0	0
2	B	214	Total 1689	C 1097	N 284	O 296	S 12	0	0	0
2	C	224	Total 1781	C 1160	N 298	O 311	S 12	0	0	0
2	D	223	Total 1769	C 1146	N 298	O 313	S 12	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	33	Total 33	O 33	0	0
3	A	97	Total 97	O 97	0	0
3	B	17	Total 17	O 17	0	0
3	C	16	Total 16	O 16	0	0

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.
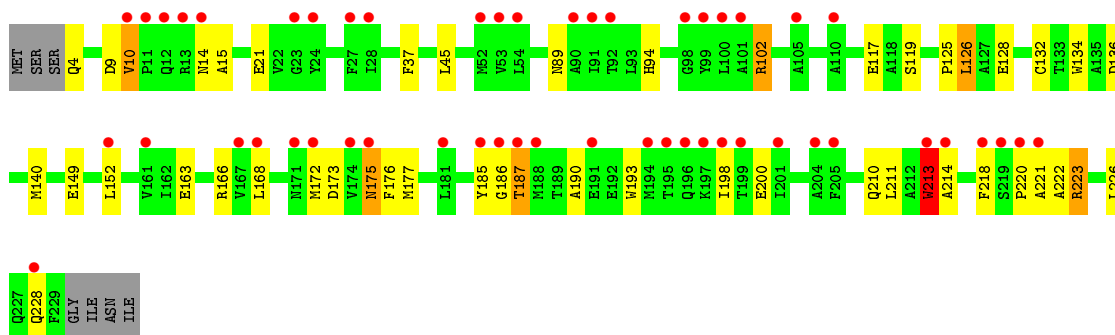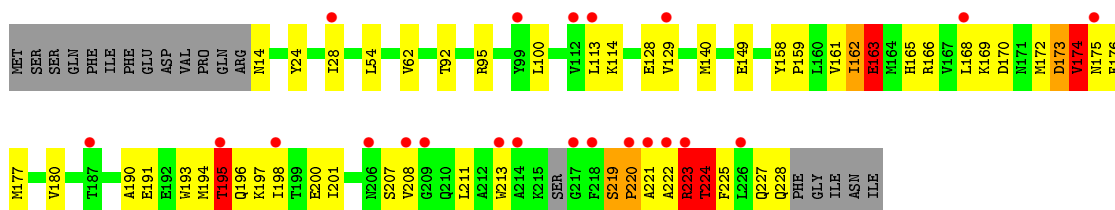
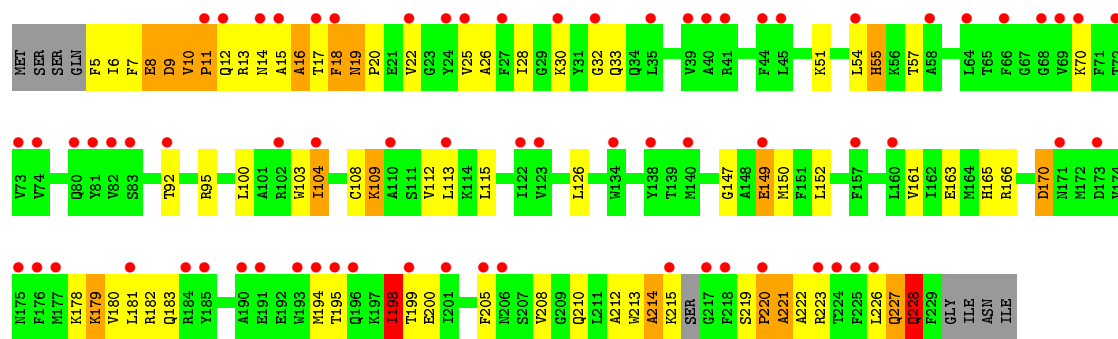- Molecule 1: RNA (42-MER)
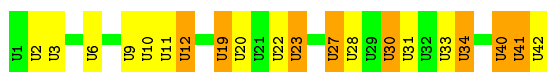
- Molecule 2: Nucleocapsid protein

- Molecule 2: Nucleocapsid protein

- Molecule 2: Nucleocapsid protein

• Molecule 2: Nucleocapsid protein

Chain D:
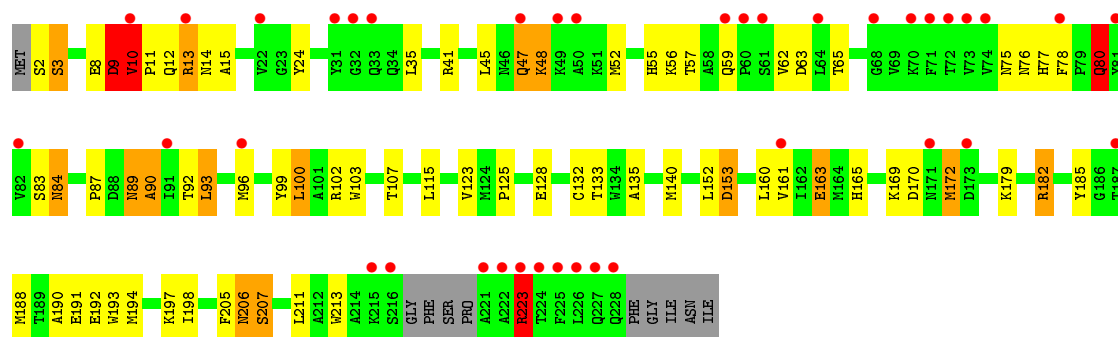
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	76.35Å 85.96Å 77.33Å 90.00° 102.02° 90.00°	Depositor
Resolution (Å)	37.40 – 2.12 37.37 – 2.12	Depositor EDS
% Data completeness (in resolution range)	97.7 (37.40-2.12) 97.7 (37.37-2.12)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	2.39 (at 2.12Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.264 , 0.283 0.280 , 0.301	Depositor DCC
R_{free} test set	2755 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	36.6	Xtriage
Anisotropy	0.653	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 37.0	EDS
L-test for twinning[2]	$< |L| > = 0.48, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.023 for l,-k,h	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	8038	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

[2]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.

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	L	0.32	0/923	0.70	0/1424
2	A	0.36	0/1844	0.64	0/2497
2	B	0.36	0/1732	0.65	1/2344 (0.0%)
2	C	0.33	0/1828	0.61	0/2474
2	D	0.35	0/1813	0.68	3/2453 (0.1%)
All	All	0.35	0/8140	0.65	4/11192 (0.0%)

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 mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	223	ARG	NE-CZ-NH1	6.02	123.31	120.30
2	D	206	ASN	N-CA-C	5.78	126.60	111.00
2	D	206	ASN	C-N-CA	5.23	134.77	121.70
2	D	10	VAL	C-N-CD	-5.11	109.36	120.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	224	THR	Peptide

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	L	840	0	421	29	0
2	A	1796	0	1781	34	0
2	B	1689	0	1682	47	0
2	C	1781	0	1767	53	0
2	D	1769	0	1757	60	0
3	A	97	0	0	3	0
3	B	17	0	0	0	0
3	C	16	0	0	0	0
3	L	33	0	0	0	0
All	All	8038	0	7408	192	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 13.

The worst 5 of 192 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:B:220:PRO:HB2	2:B:221:ALA:HA	1.28	1.15
2:C:12:GLN:HB3	2:C:13:ARG:HA	1.11	1.10
2:A:9:ASP:HA	2:A:10:VAL:HG22	1.36	1.07
1:L:12:U:H5'	1:L:12:U:H6	1.24	1.03
2:B:162:ILE:O	2:B:163:GLU:HB2	1.55	1.01

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	
2	A	224/233 (96%)	198 (88%)	19 (8%)	7 (3%)	4	1
2	B	210/233 (90%)	181 (86%)	21 (10%)	8 (4%)	3	1
2	C	220/233 (94%)	184 (84%)	19 (9%)	17 (8%)	1	0
2	D	219/233 (94%)	190 (87%)	15 (7%)	14 (6%)	1	0
All	All	873/932 (94%)	753 (86%)	74 (8%)	46 (5%)	2	0

5 of 46 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	126	LEU
2	A	213	TRP
2	A	221	ALA
2	B	163	GLU
2	B	174	VAL

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	
2	A	187/193 (97%)	173 (92%)	14 (8%)	13	10
2	B	175/193 (91%)	164 (94%)	11 (6%)	18	14
2	C	185/193 (96%)	163 (88%)	22 (12%)	5	2
2	D	185/193 (96%)	165 (89%)	20 (11%)	6	3
All	All	732/772 (95%)	665 (91%)	67 (9%)	9	5

5 of 67 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	57	THR
2	C	163	GLU
2	D	170	ASP
2	C	70	LYS
2	C	126	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	94	HIS
2	B	206	ASN
2	D	47	GLN
2	B	59	GLN
2	D	55	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	L	41/42 (97%)	11 (26%)	6 (14%)

5 of 11 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	L	6	U
1	L	11	U
1	L	12	U
1	L	19	U
1	L	22	U

5 of 6 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	L	22	U
1	L	40	U
1	L	23	U
1	L	12	U
1	L	33	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 (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, 95th 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	L	42/42 (100%)	-0.54	0	100	100	5, 18, 32, 54	0
2	A	226/233 (96%)	1.31	51 (22%)	0	0	5, 10, 40, 49	0
2	B	214/233 (91%)	1.04	22 (10%)	6	8	5, 13, 42, 67	0
2	C	224/233 (96%)	1.73	71 (31%)	0	0	10, 28, 49, 78	0
2	D	223/233 (95%)	1.16	38 (17%)	1	1	5, 16, 41, 52	0
All	All	929/974 (95%)	1.23	182 (19%)	1	1	5, 17, 43, 78	0

The worst 5 of 182 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	221	ALA	15.1
2	C	218	PHE	9.3
2	C	191	GLU	8.6
2	B	213	TRP	8.5
2	A	167	VAL	6.6

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.