



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 12:23 pm BST

PDB ID : 5JPS
Title : Structure of the Zika Virus NS3 Helicase Domain
Authors : Jia, Z.; Wang, J.
Deposited on : 2016-05-04
Resolution : 1.78 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 ⓘ 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

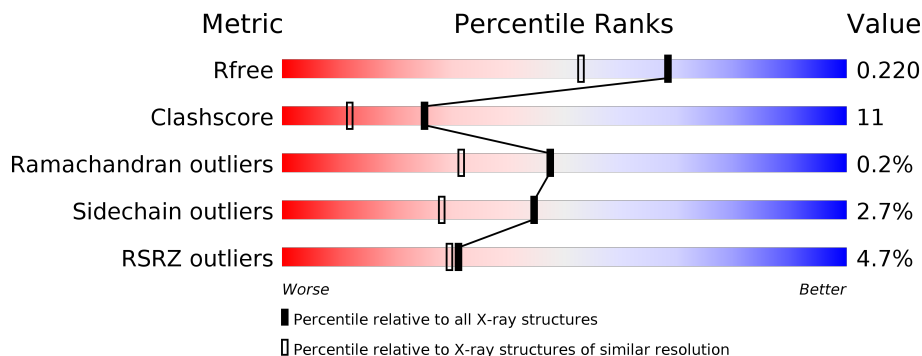
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

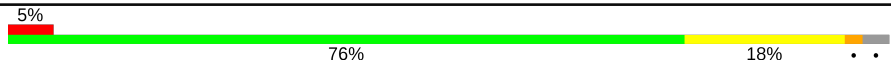
The reported resolution of this entry is 1.78 Å.

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}	130704	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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 $\leq 5\%$. 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	442	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3804 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 ns3 helicase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	430	3406	2145	608	631	22	0	0	0

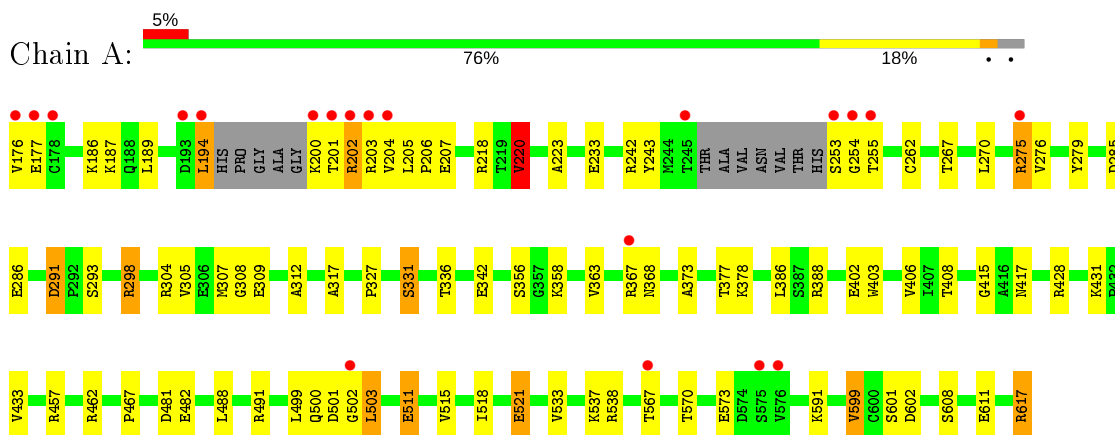
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	398	Total	O	0	0
			398	398		

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: ns3 helicase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.55Å 69.04Å 57.15Å 90.00° 91.91° 90.00°	Depositor
Resolution (Å)	42.33 – 1.78 39.72 – 1.78	Depositor EDS
% Data completeness (in resolution range)	98.5 (42.33-1.78) 98.5 (39.72-1.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 1.78Å)	Xtrriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.180 , 0.215 0.188 , 0.220	Depositor DCC
R_{free} test set	1963 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	18.9	Xtrriage
Anisotropy	0.078	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.030 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3804	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	1.02	6/3476 (0.2%)	1.00	15/4701 (0.3%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	298	ARG	CD-NE	-11.88	1.26	1.46
1	A	327	PRO	C-N	9.57	1.56	1.34
1	A	617	ARG	CZ-NH2	-5.97	1.25	1.33
1	A	331	SER	CB-OG	-5.74	1.34	1.42
1	A	457	ARG	CZ-NH2	5.23	1.39	1.33
1	A	511	GLU	CB-CG	-5.11	1.42	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	298	ARG	NE-CZ-NH2	-23.48	108.56	120.30
1	A	298	ARG	NE-CZ-NH1	14.60	127.60	120.30
1	A	617	ARG	NE-CZ-NH2	-10.96	114.82	120.30
1	A	298	ARG	CD-NE-CZ	9.75	137.25	123.60
1	A	218	ARG	NE-CZ-NH2	-7.74	116.43	120.30
1	A	218	ARG	NE-CZ-NH1	7.27	123.93	120.30
1	A	599	VAL	CG1-CB-CG2	7.18	122.39	110.90
1	A	298	ARG	CG-CD-NE	-7.04	97.00	111.80
1	A	617	ARG	NE-CZ-NH1	6.99	123.80	120.30
1	A	491	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	A	220	VAL	CG1-CB-CG2	6.25	120.90	110.90
1	A	291	ASP	CB-CG-OD2	-6.25	112.68	118.30
1	A	220	VAL	CB-CA-C	-6.21	99.61	111.40
1	A	428	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	A	457	ARG	NE-CZ-NH2	-5.32	117.64	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	3406	0	3402	73	0
2	A	398	0	0	21	3
All	All	3804	0	3402	73	3

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 11.

All (73) 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:A:406:VAL:HG12	1:A:408:THR:HG23	1.60	0.83
1:A:200:LYS:N	1:A:202:ARG:HG3	1.96	0.81
1:A:573:GLU:OE1	2:A:701:HOH:O	2.01	0.78
1:A:275:ARG:HG2	1:A:275:ARG:HH11	1.51	0.75
1:A:275:ARG:NE	2:A:706:HOH:O	2.18	0.75
1:A:189:LEU:HD13	1:A:312:ALA:HB3	1.70	0.74
1:A:307:MET:SD	1:A:499:LEU:HD21	2.28	0.73
1:A:186:LYS:O	1:A:187:LYS:HB2	1.88	0.73
1:A:200:LYS:N	1:A:202:ARG:CG	2.51	0.72
1:A:267:THR:HG21	1:A:293:SER:O	1.93	0.69
1:A:298:ARG:HD2	1:A:511:GLU:OE2	1.93	0.69
1:A:482:GLU:OE1	2:A:702:HOH:O	2.10	0.69
1:A:367:ARG:HE	1:A:602:ASP:HA	1.58	0.69
1:A:336:THR:HG23	2:A:803:HOH:O	1.92	0.68
1:A:253:SER:N	2:A:712:HOH:O	2.26	0.68
1:A:200:LYS:HE3	1:A:285:ASP:OD1	1.95	0.66
1:A:402:GLU:OE2	2:A:703:HOH:O	2.12	0.66
1:A:481:ASP:HB2	2:A:918:HOH:O	1.95	0.66
1:A:187:LYS:NZ	2:A:716:HOH:O	2.30	0.64
1:A:567:THR:O	1:A:570:THR:HG23	1.99	0.63
1:A:367:ARG:NE	1:A:601:SER:O	2.33	0.62
1:A:406:VAL:HG12	1:A:408:THR:CG2	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:VAL:HG22	1:A:279:TYR:CG	2.37	0.60
1:A:481:ASP:OD2	2:A:705:HOH:O	2.16	0.59
1:A:242:ARG:NH1	2:A:709:HOH:O	2.20	0.59
1:A:275:ARG:CD	2:A:706:HOH:O	2.51	0.59
1:A:275:ARG:HD3	2:A:706:HOH:O	2.05	0.57
1:A:275:ARG:CZ	2:A:706:HOH:O	2.53	0.57
1:A:611:GLU:OE1	2:A:707:HOH:O	2.18	0.56
1:A:533:VAL:O	1:A:537:LYS:HG2	2.05	0.56
1:A:417:ASN:OD1	1:A:462:ARG:NH1	2.35	0.56
1:A:608:SER:O	1:A:617:ARG:NH2	2.33	0.56
1:A:176:VAL:HG13	1:A:177:GLU:H	1.71	0.55
1:A:342:GLU:OE1	2:A:708:HOH:O	2.18	0.54
1:A:386:LEU:HB2	1:A:408:THR:HG22	1.89	0.53
1:A:431:LYS:HE3	1:A:433:VAL:HG22	1.91	0.53
1:A:187:LYS:HG2	1:A:308:GLY:HA2	1.91	0.52
1:A:220:VAL:HG22	1:A:279:TYR:CD1	2.45	0.52
1:A:502:GLY:C	1:A:503:LEU:HD23	2.29	0.52
1:A:358:LYS:HE2	1:A:403:TRP:CE2	2.46	0.51
1:A:538:ARG:HG3	1:A:538:ARG:HH11	1.76	0.51
1:A:518:ILE:O	1:A:521:GLU:HB3	2.12	0.50
1:A:500:GLN:O	1:A:501:ASP:C	2.48	0.50
1:A:233:GLU:HG2	1:A:243:TYR:CE1	2.48	0.49
1:A:255:THR:HA	2:A:710:HOH:O	2.11	0.49
1:A:373:ALA:O	1:A:377:THR:HG23	2.12	0.49
1:A:255:THR:OG1	1:A:255:THR:O	2.28	0.49
1:A:462:ARG:NH2	2:A:704:HOH:O	2.15	0.48
1:A:367:ARG:CZ	1:A:601:SER:O	2.61	0.48
1:A:488:LEU:HA	1:A:515:VAL:HG21	1.96	0.48
1:A:304:ARG:NH1	1:A:309:GLU:OE1	2.45	0.48
1:A:406:VAL:CG1	1:A:408:THR:CG2	2.93	0.47
1:A:203:ARG:O	1:A:207:GLU:HG3	2.15	0.47
1:A:205:LEU:HB3	1:A:206:PRO:HD3	1.95	0.47
1:A:275:ARG:CG	1:A:275:ARG:HH11	2.25	0.46
1:A:200:LYS:HB3	1:A:201:THR:H	1.40	0.45
1:A:194:LEU:HB3	1:A:317:ALA:CB	2.48	0.44
1:A:223:ALA:O	1:A:262:CYS:HA	2.17	0.44
1:A:200:LYS:HE2	1:A:286:GLU:OE2	2.17	0.43
1:A:431:LYS:NZ	1:A:482:GLU:O	2.51	0.43
1:A:270:LEU:CD2	1:A:276:VAL:HG21	2.49	0.43
1:A:358:LYS:HE2	1:A:403:TRP:CD2	2.54	0.43
1:A:570:THR:HG22	1:A:591:LYS:HD3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:LEU:HD23	1:A:276:VAL:HG21	2.00	0.42
1:A:200:LYS:N	1:A:202:ARG:HG2	2.32	0.41
1:A:363:VAL:HB	1:A:368:ASN:HD22	1.85	0.41
1:A:342:GLU:HG2	2:A:732:HOH:O	2.20	0.41
1:A:331:SER:OG	1:A:467:PRO:HA	2.21	0.41
1:A:202:ARG:HB2	1:A:204:VAL:H	1.85	0.41
1:A:415:GLY:HA2	2:A:704:HOH:O	2.21	0.41
1:A:254:GLY:O	2:A:710:HOH:O	2.22	0.40
1:A:203:ARG:C	1:A:206:PRO:HD2	2.42	0.40
1:A:356:SER:O	2:A:711:HOH:O	2.22	0.40

All (3) 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:A:897:HOH:O	2:A:1055:HOH:O[1_556]	2.03	0.17
2:A:703:HOH:O	2:A:992:HOH:O[1_554]	2.14	0.06
2:A:1046:HOH:O	2:A:1079:HOH:O[1_554]	2.14	0.06

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	424/442 (96%)	411 (97%)	12 (3%)	1 (0%)	47 32

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	202	ARG

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	364/372 (98%)	354 (97%)	10 (3%)	44 28

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	194	LEU
1	A	220	VAL
1	A	275	ARG
1	A	291	ASP
1	A	305	VAL
1	A	378	LYS
1	A	388	ARG
1	A	503	LEU
1	A	521	GLU
1	A	599	VAL

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	368	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 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	A	430/442 (97%)	0.21	20 (4%) 31 29	10, 19, 35, 58	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	176	VAL	7.1
1	A	203	ARG	5.8
1	A	253	SER	5.1
1	A	275	ARG	5.0
1	A	194	LEU	5.0
1	A	201	THR	4.0
1	A	178	CYS	3.8
1	A	254	GLY	3.8
1	A	255	THR	3.8
1	A	502	GLY	3.7
1	A	177	GLU	3.3
1	A	200	LYS	3.2
1	A	193	ASP	3.1
1	A	576	VAL	2.9
1	A	367	ARG	2.7
1	A	567	THR	2.7
1	A	575	SER	2.2
1	A	204	VAL	2.1
1	A	202	ARG	2.0
1	A	245	THR	2.0

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.