

wwPDB X-ray Structure Validation Summary Report (i)

Oct 31, 2023 – 03:01 PM JST

PDB ID : 5GJ4

Title : Structure of NS2B-NS3 Protease from Zika Virus caught after self-cleavage

Authors : Phoo, W.W.; Li, Y.

Deposited on : 2016-06-27

Resolution : 1.84 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 (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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

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

EDS : 2.36

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)

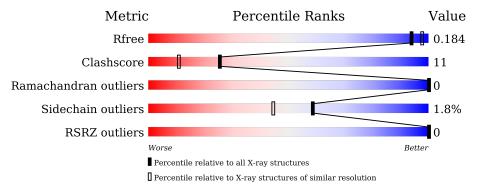
Validation Pipeline (wwPDB-VP) : 2.36

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 1.84 Å.

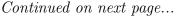
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}({\rm \AA})) \end{array}$
R_{free}	130704	4003 (1.86-1.82)
Clashscore	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)
RSRZ outliers	127900	3957 (1.86-1.82)

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 >=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	A	61	61%	10%	30%				
1	С	61	61%	8% •	30%				
1	Е	61	51%	18% •	30%				
1	G	61	56%	15%	30%				
2	В	177	73%		14% 14%				
2	D	177	71%	•	15% • 14%				





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Mol	Chain	Length	Quality of chain					
2	F	177	72%	14%	·	14%		
2	Н	177	71%	15%		14%		



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7043 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 Serine protease subunit NS2B.

Mol	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
1	Λ	43	Total	С	N	О	S	0	0	0
1	A	40	334	205	56	72	1	U		
1	С	43	Total	С	N	О	S	0	0	0
1		40	334	205	56	72	1			
1	Е	43	Total	С	N	О	S	0	0	0
1	12	40	334	205	56	72	1	0		
1	G	43	Total	С	N	О	S	0	0	0
1	G	40	334	205	56	72	1	U		U

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	GLY	-	expression tag	UNP A0A142IX72
A	42	SER	-	expression tag	UNP A0A142IX72
A	43	HIS	-	expression tag	UNP A0A142IX72
A	44	MET	-	expression tag	UNP A0A142IX72
С	41	GLY	_	expression tag	UNP A0A142IX72
С	42	SER	-	expression tag	UNP A0A142IX72
С	43	HIS	-	expression tag	UNP A0A142IX72
С	44	MET	-	expression tag	UNP A0A142IX72
Е	41	GLY	-	expression tag	UNP A0A142IX72
Е	42	SER	_	expression tag	UNP A0A142IX72
Е	43	HIS	-	expression tag	UNP A0A142IX72
Е	44	MET	-	expression tag	UNP A0A142IX72
G	41	GLY	-	expression tag	UNP A0A142IX72
G	42	SER	-	expression tag	UNP A0A142IX72
G	43	HIS	-	expression tag	UNP A0A142IX72
G	44	MET	_	expression tag	UNP A0A142IX72

• Molecule 2 is a protein called Serine protease NS3.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	В	153	Total	С	N	О	S	0	0	0
2	Б	199	1144	724	202	213	5	0	U	U
2	D	152	Total	С	N	О	S	0	0	0
	D	192	1139	721	201	212	5			
2	F	152	Total	С	N	О	S	0	1	0
	I'	192	1140	721	200	214	5	0		U
2	Н	153	Total	С	N	О	S	0	0	0
2	11	199	1144	724	202	213	5			U

• Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	В	1	Total Cl 1 1	0	0
3	D	1	Total Cl 1 1	0	0

• Molecule 4 is water.

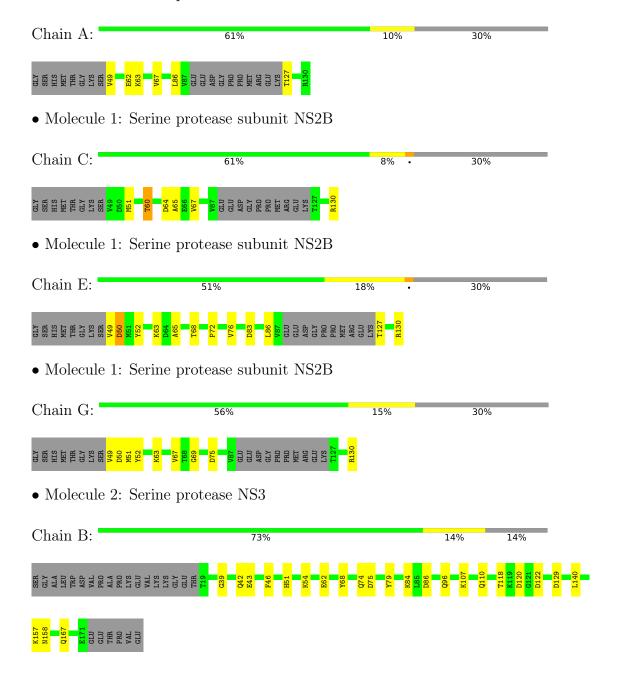
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	89	Total O 89 89	0	0
4	В	199	Total O 199 199	0	0
4	С	81	Total O 81 81	0	0
4	D	199	Total O 199 199	0	0
4	E	101	Total O 101 101	0	0
4	F	210	Total O 210 210	0	0
4	G	75	Total O 75 75	0	0
4	Н	184	Total O 184 184	0	0



3 Residue-property plots (i)

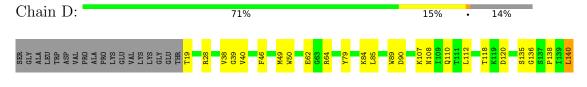
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Serine protease subunit NS2B



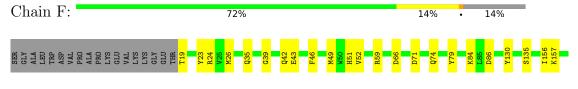


• Molecule 2: Serine protease NS3



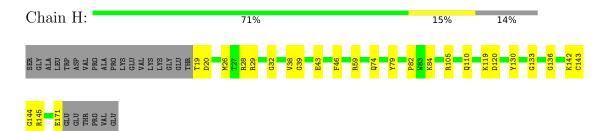


• Molecule 2: Serine protease NS3





• Molecule 2: Serine protease NS3





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants	88.79Å 88.79Å 138.10Å	Donositon
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	31.50 - 1.84	Depositor
rtesolution (A)	31.50 - 1.84	EDS
% Data completeness	97.8 (31.50-1.84)	Depositor
(in resolution range)	94.6 (31.50-1.84)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.26 (at 1.84Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.169 , 0.195	Depositor
	0.161 , 0.184	DCC
R_{free} test set	5122 reflections $(4.84%)$	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.28, 32.1	EDS
L-test for twinning ²	$< L > = 0.39, < L^2> = 0.22$	Xtriage
	0.426 for -h,-k,l	
Estimated twinning fraction	0.487 for h,-h-k,-l	Xtriage
	0.427 for -k,-h,-l	
F_o, F_c correlation	0.96	EDS
Total number of atoms	7043	wwPDB-VP
Average B, all atoms (\mathring{A}^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CL

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	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.30	0/337	0.53	0/452	
1	С	0.30	0/337	0.51	0/452	
1	Е	0.30	0/337	0.52	0/452	
1	G	0.30	0/337	0.52	0/452	
2	В	0.31	0/1168	0.54	0/1585	
2	D	0.31	0/1163	0.56	1/1578~(0.1%)	
2	F	0.30	0/1167	0.54	0/1585	
2	Н	0.30	0/1168	0.54	0/1585	
All	All	0.31	0/6014	0.54	1/8141 (0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^o)$
2	D	140	LEU	CA-CB-CG	5.33	127.55	115.30

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	334	0	315	9	0
1	С	334	0	315	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Ε	334	0	315	15	0
1	G	334	0	315	16	0
2	В	1144	0	1139	21	0
2	D	1139	0	1137	27	0
2	F	1140	0	1130	23	0
2	Η	1144	0	1139	30	0
3	В	1	0	0	0	0
3	D	1	0	0	0	0
4	A	89	0	0	6	0
4	В	199	0	0	12	3
4	С	81	0	0	4	0
4	D	199	0	0	9	3
4	Ε	101	0	0	8	1
4	F	210	0	0	8	2
4	G	75	0	0	9	0
4	Н	184	0	0	10	3
All	All	7043	0	5805	124	6

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.

The worst 5 of 124 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
2:F:86:ASP:OD1	4:F:201:HOH:O	1.85	0.94
1:A:49:VAL:O	4:A:201:HOH:O	1.96	0.82
2:B:107:LYS:NZ	4:B:2105:HOH:O	2.12	0.81
1:A:49:VAL:N	4:A:202:HOH:O	2.13	0.80
2:B:167:GLN:O	4:B:2101:HOH:O	1.98	0.80

The worst 5 of 6 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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
4:B:2204:HOH:O	4:F:366:HOH:O[3_554]	1.80	0.40
4:B:2204:HOH:O	4:F:344:HOH:O[3_554]	1.88	0.32
4:D:336:HOH:O	4:H:251:HOH:O[2_665]	2.09	0.11
4:B:2242:HOH:O	4:H:370:HOH:O[3_564]	2.12	0.08
4:D:485:HOH:O	4:H:384:HOH:O[2_665]	2.15	0.05



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	Perce	entiles
1	A	39/61 (64%)	39 (100%)	0	0	100	100
1	С	39/61 (64%)	38 (97%)	1 (3%)	0	100	100
1	E	39/61 (64%)	38 (97%)	1 (3%)	0	100	100
1	G	39/61 (64%)	39 (100%)	0	0	100	100
2	В	151/177 (85%)	147 (97%)	4 (3%)	0	100	100
2	D	150/177~(85%)	146 (97%)	4 (3%)	0	100	100
2	F	151/177 (85%)	145 (96%)	6 (4%)	0	100	100
2	Н	151/177 (85%)	146 (97%)	5 (3%)	0	100	100
All	All	759/952 (80%)	738 (97%)	21 (3%)	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	lysed Rotameric		Perce	ntiles
1	A	36/51 (71%)	36 (100%)	0	100	100
1	С	36/51 (71%)	34 (94%)	2 (6%)	21	6
1	E	36/51 (71%)	34 (94%)	2 (6%)	21	6
1	G	36/51 (71%)	35 (97%)	1 (3%)	43	26
2	В	118/140 (84%)	116 (98%)	2 (2%)	60	47
2	D	118/140 (84%)	117 (99%)	1 (1%)	81	75

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
2	F	118/140 (84%)	117 (99%)	1 (1%)	81 75		
2	Н	118/140 (84%)	116 (98%)	2 (2%)	60 47		
All	All	616/764 (81%)	605 (98%)	11 (2%)	59 44		

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	F	170	ARG
1	G	51	MET
2	Н	120	ASP
2	Н	119	LYS
2	D	140	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	В	158	ASN
2	D	108	ASN
2	F	35	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.



There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	<RSRZ $>$	$\#\mathrm{RSRZ}{>}2$		ZZ>2	$OWAB(A^2)$	Q < 0.9
1	A	43/61 (70%)	-0.51	0	100	100	17, 32, 40, 45	0
1	С	43/61 (70%)	-0.55	0	100	100	18, 33, 42, 47	0
1	E	43/61 (70%)	-0.43	0	100	100	17, 32, 45, 53	0
1	G	43/61 (70%)	-0.42	0	100	100	18, 33, 43, 52	0
2	В	153/177 (86%)	-0.61	0	100	100	16, 25, 40, 59	0
2	D	152/177~(85%)	-0.62	0	100	100	16, 25, 36, 42	0
2	F	152/177~(85%)	-0.63	0	100	100	15, 25, 38, 51	0
2	Н	153/177 (86%)	-0.64	0	100	100	17, 24, 37, 58	0
All	All	782/952~(82%)	-0.59	0	100	100	15, 26, 41, 59	0

There are no RSRZ outliers to report.

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 monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.



Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q<0.9
3	CL	В	2001	1/1	0.99	0.07	19,19,19,19	0
3	CL	D	201	1/1	0.99	0.07	19,19,19,19	0

6.5 Other polymers (i)

There are no such residues in this entry.

