



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:39 PM GMT

PDB ID : 5A9B  
Title : Crystal structure of Bombyx mori CPV1 polyhedra base domain deleted mutant  
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Deposited on : 2015-07-17  
Resolution : 1.88 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](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  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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-20026982

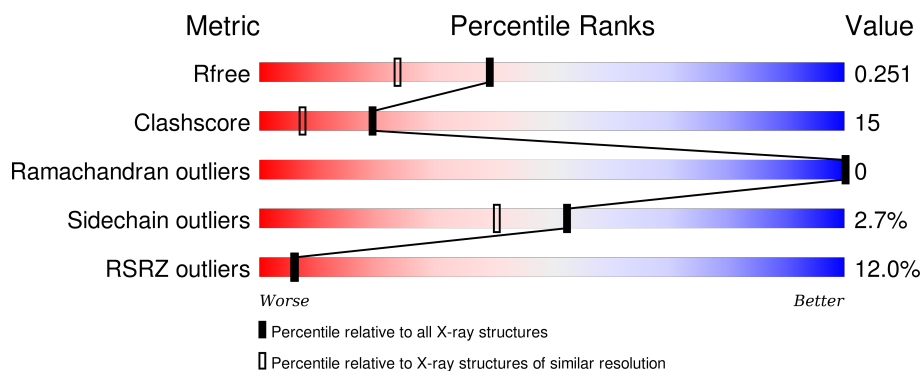
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.88 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))
$R_{free}$	91344	6965 (1.90-1.86)
Clashscore	102246	7778 (1.90-1.86)
Ramachandran outliers	100387	7691 (1.90-1.86)
Sidechain outliers	100360	7692 (1.90-1.86)
RSRZ outliers	91569	6979 (1.90-1.86)

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	218	<div> <div>11%</div> <div>74%</div> <div>21%</div> <div>..</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1787 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 POLYHEDRIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	1	0
			1686	1069	295	318	4			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	SER	ASN	CONFLICT	UNP P11041
A	67	VAL	ALA	CONFLICT	UNP P11041
A	.	-	TYR	DELETION	UNP P11041
A	.	-	ARG	DELETION	UNP P11041
A	.	-	GLU	DELETION	UNP P11041
A	.	-	GLY	DELETION	UNP P11041
A	.	-	GLN	DELETION	UNP P11041
A	.	-	HIS	DELETION	UNP P11041
A	.	-	ASN	DELETION	UNP P11041
A	.	-	ASP	DELETION	UNP P11041
A	.	-	SER	DELETION	UNP P11041
A	.	-	TYR	DELETION	UNP P11041
A	.	-	ASP	DELETION	UNP P11041
A	.	-	GLU	DELETION	UNP P11041
A	.	-	TYR	DELETION	UNP P11041
A	.	-	GLU	DELETION	UNP P11041
A	.	-	VAL	DELETION	UNP P11041
A	.	-	ASN	DELETION	UNP P11041
A	.	-	GLN	DELETION	UNP P11041
A	.	-	SER	DELETION	UNP P11041
A	.	-	ILE	DELETION	UNP P11041
A	.	-	TYR	DELETION	UNP P11041
A	.	-	TYR	DELETION	UNP P11041
A	.	-	PRO	DELETION	UNP P11041
A	.	-	ASN	DELETION	UNP P11041
A	.	-	GLY	DELETION	UNP P11041
A	.	-	GLY	DELETION	UNP P11041

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Chain	Residue	Modelled	Actual	Comment	Reference
A	.	-	ASP	DELETION	UNP P11041
A	.	-	ALA	DELETION	UNP P11041
A	.	-	ARG	DELETION	UNP P11041
A	.	-	LYS	DELETION	UNP P11041
A	.	-	PHE	DELETION	UNP P11041
A	.	-	HIS	DELETION	UNP P11041
A	.	-	SER	DELETION	UNP P11041
A	.	-	ASN	DELETION	UNP P11041
A	101	GLY	-	INSERTION	UNP P11041
A	102	SER	-	INSERTION	UNP P11041
A	103	GLY	-	INSERTION	UNP P11041

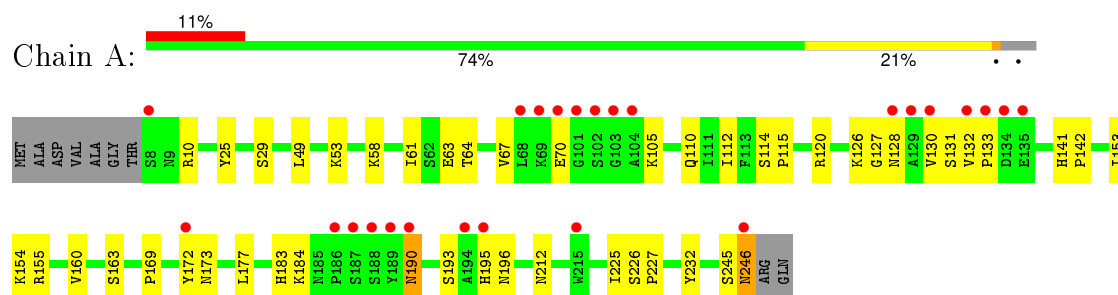
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	101	Total O 101 101	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 errors 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 ( $\text{RSRZ} > 2$ ). 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: POLYHEDRIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.86Å 103.86Å 103.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.40 – 1.88 42.40 – 1.88	Depositor EDS
% Data completeness (in resolution range)	98.4 (42.40-1.88) 98.4 (42.40-1.88)	Depositor EDS
$R_{merge}$	0.23	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.22 (at 1.88Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.189 , 0.251 0.187 , 0.251	Depositor DCC
$R_{free}$ test set	758 reflections (5.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 64.0	EDS
Estimated twinning fraction	0.060 for -l,-k,-h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 15018 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	1787	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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	0.26	0/1733	0.44	0/2350

There are no bond length outliers.

There are no bond angle outliers.

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	1686	0	1622	50	0
2	A	101	0	0	5	1
All	All	1787	0	1622	50	1

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 15.

All (50) 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:127:GLY:CA	1:A:133:PRO:HG3	1.49	1.42
1:A:127:GLY:CA	1:A:133:PRO:CG	2.12	1.28
1:A:127:GLY:HA3	1:A:133:PRO:CG	1.77	1.14
1:A:127:GLY:HA3	1:A:133:PRO:HG3	1.25	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:GLY:HA2	1:A:133:PRO:CG	1.89	1.02
1:A:127:GLY:HA2	1:A:133:PRO:HG2	1.42	1.01
1:A:127:GLY:C	1:A:133:PRO:HG3	1.81	0.98
1:A:141:HIS:CD2	1:A:142:PRO:HD2	1.98	0.98
1:A:127:GLY:CA	1:A:133:PRO:HG2	1.99	0.89
1:A:131:SER:O	1:A:132:VAL:CG2	2.30	0.81
1:A:131:SER:O	1:A:132:VAL:HG23	1.82	0.78
1:A:67:VAL:HG23	1:A:173:ASN:O	1.85	0.77
1:A:67:VAL:HG22	1:A:105:LYS:HB3	1.66	0.76
1:A:193:SER:OG	1:A:196:ASN:ND2	2.20	0.75
1:A:128:ASN:N	1:A:133:PRO:HG3	2.05	0.71
1:A:53:LYS:HB2	1:A:61:ILE:HD11	1.71	0.71
1:A:131:SER:C	1:A:132:VAL:HG22	2.11	0.71
1:A:173:ASN:ND2	2:A:2041:HOH:O	2.21	0.69
1:A:25:TYR:OH	2:A:2010:HOH:O	2.11	0.68
1:A:131:SER:C	1:A:132:VAL:CG2	2.64	0.66
1:A:128:ASN:H	1:A:133:PRO:CB	2.11	0.62
1:A:190:ASN:HB3	1:A:193:SER:HB3	1.81	0.60
1:A:183:HIS:ND1	1:A:184:LYS:O	2.32	0.60
1:A:169:PRO:HB2	1:A:172:TYR:HD2	1.68	0.59
1:A:64:THR:OG1	2:A:2039:HOH:O	2.17	0.58
1:A:110:GLN:OE1	1:A:112:ILE:HD11	2.05	0.57
1:A:67:VAL:HG22	1:A:105:LYS:CB	2.33	0.56
1:A:61:ILE:CG2	1:A:63:GLU:HG2	2.36	0.56
1:A:126:LYS:HB3	1:A:177:LEU:HD12	1.89	0.54
1:A:128:ASN:N	1:A:133:PRO:CG	2.70	0.54
1:A:225:ILE:HD12	1:A:232:TYR:HB2	1.93	0.51
1:A:169:PRO:HB2	1:A:172:TYR:CD2	2.46	0.51
1:A:70:GLU:HG3	2:A:2042:HOH:O	2.11	0.50
1:A:153:ILE:HD11	1:A:163:SER:HB2	1.93	0.49
1:A:128:ASN:H	1:A:133:PRO:HB2	1.78	0.48
1:A:127:GLY:HA3	1:A:133:PRO:CD	2.40	0.48
1:A:130:VAL:HG12	1:A:130:VAL:O	2.13	0.47
1:A:128:ASN:N	1:A:133:PRO:CB	2.78	0.47
1:A:53:LYS:HE3	1:A:61:ILE:HD13	1.97	0.47
1:A:49:LEU:HD12	1:A:212:ASN:O	2.15	0.47
1:A:10:ARG:HH11	1:A:10:ARG:HG3	1.80	0.46
1:A:245:SER:HA	1:A:246:ASN:HA	1.61	0.46
1:A:226:SER:HB2	1:A:227:PRO:HD3	1.99	0.45
1:A:29:SER:CB	2:A:2016:HOH:O	2.57	0.45
1:A:61:ILE:HG23	1:A:63:GLU:HG2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:LYS:O	1:A:155:ARG:HD2	2.18	0.44
1:A:128:ASN:H	1:A:133:PRO:CG	2.32	0.43
1:A:115:PRO:HD3	1:A:160:VAL:O	2.20	0.41
1:A:114:SER:HB2	1:A:115:PRO:HD2	2.02	0.41
1:A:131:SER:O	1:A:132:VAL:HG22	2.09	0.40

All (1) 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:2059:HOH:O	2:A:2095:HOH:O[24_554]	2.02	0.18

## 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	208/218 (95%)	201 (97%)	7 (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	Rotameric	Outliers	Percentiles	
1	A	185/190 (97%)	180 (97%)	5 (3%)	52	40

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	LYS
1	A	120	ARG
1	A	190	ASN
1	A	195	HIS
1	A	246	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

### 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	209/218 (95%)	0.38	25 (11%) 6 6	16, 29, 80, 105	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	133	PRO	7.9
1	A	132	VAL	7.4
1	A	186	PRO	7.0
1	A	129	ALA	6.9
1	A	134	ASP	6.7
1	A	130	VAL	6.1
1	A	101	GLY	5.2
1	A	70	GLU	4.6
1	A	102	SER	4.6
1	A	68	LEU	4.4
1	A	189	TYR	4.3
1	A	187	SER	4.0
1	A	172	TYR	3.8
1	A	69	LYS	3.8
1	A	8	SER	3.5
1	A	195	HIS	3.3
1	A	188	SER	3.3
1	A	215	TRP	2.8
1	A	246	ASN	2.7
1	A	103	GLY	2.4
1	A	190	ASN	2.4
1	A	194	ALA	2.2
1	A	135	GLU	2.1
1	A	128	ASN	2.1
1	A	104	ALA	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.