



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 10:00 am GMT

PDB ID : 3D9X  
Title : Structure of the head of the Bartonella adhesin BadA  
Authors : Zeth, K.; Lupas, A.; Martin, J.  
Deposited on : 2008-05-28  
Resolution : 1.13 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  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
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) : recalc28949

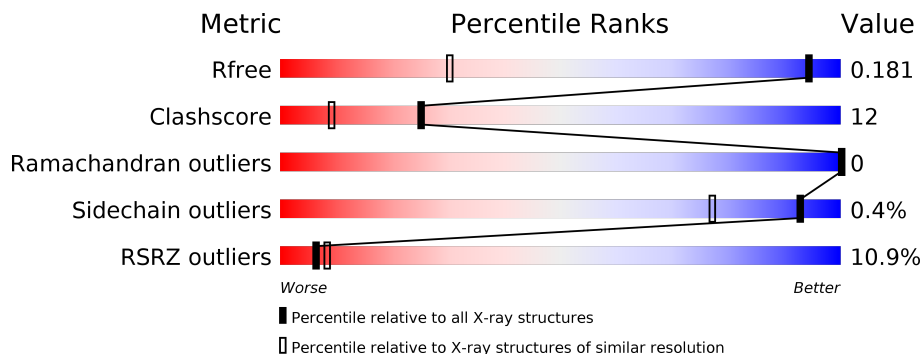
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.13 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	1062 (1.16-1.08)
Clashscore	112137	1107 (1.16-1.08)
Ramachandran outliers	110173	1065 (1.16-1.08)
Sidechain outliers	110143	1062 (1.16-1.08)
RSRZ outliers	101464	1067 (1.16-1.08)

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	163	<div> <div>8%</div> <div>65%</div> <div>5%</div> <div>30%</div> </div>
1	B	163	<div> <div>7%</div> <div>55%</div> <div>13%</div> <div>32%</div> </div>
1	C	163	<div> <div>7%</div> <div>61%</div> <div>9%</div> <div>30%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3006 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 Adhesin A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	114	Total	C	N	O	0	4	0
			853	528	149	176			
1	B	111	Total	C	N	O	0	8	0
			843	526	144	173			
1	C	114	Total	C	N	O	0	12	0
			888	559	150	179			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	374	MET	-	INITIATING METHIONINE	UNP Q5MWV9
B	374	MET	-	INITIATING METHIONINE	UNP Q5MWV9
C	374	MET	-	INITIATING METHIONINE	UNP Q5MWV9

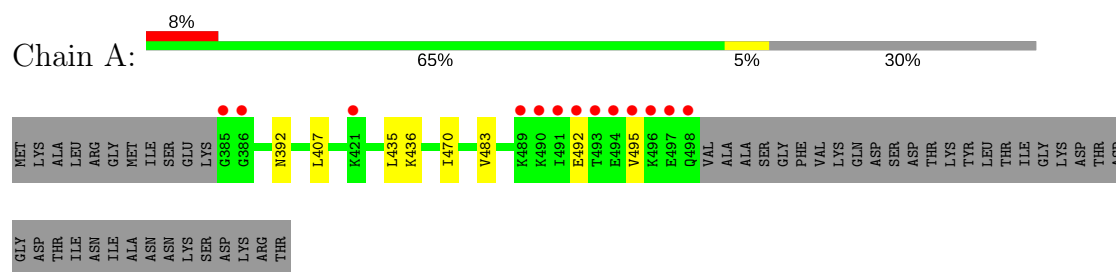
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	133	Total	O	0	0
			133	133		
2	B	133	Total	O	0	0
			133	133		
2	C	156	Total	O	0	0
			156	156		

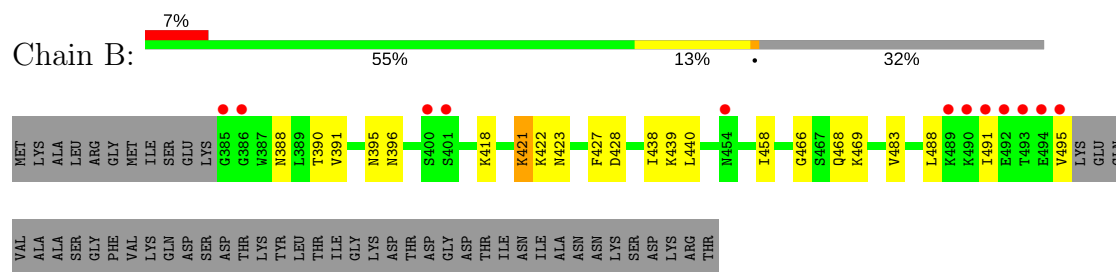
### 3 Residue-property plots [i](#)

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.

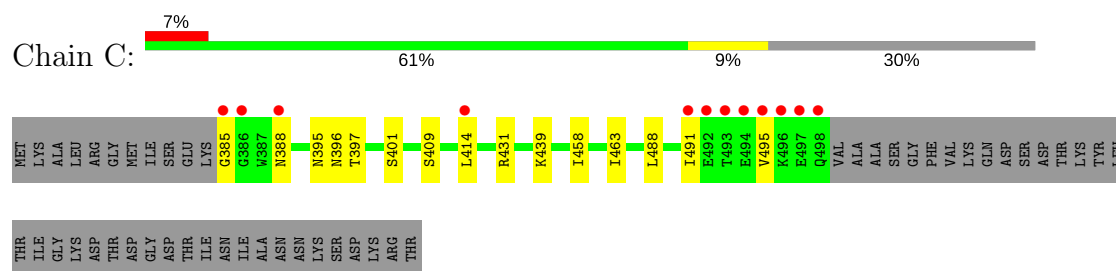
#### • Molecule 1: Adhesin A



#### • Molecule 1: Adhesin A



#### • Molecule 1: Adhesin A



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	29.87Å 51.14Å 58.62Å 65.87° 76.60° 82.08°	Depositor
Resolution (Å)	25.00 – 1.13 29.58 – 1.13	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.00-1.13) 81.8 (29.58-1.13)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.62 (at 1.13Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.156 , 0.178 0.161 , 0.181	Depositor DCC
$R_{free}$ test set	5257 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	9.4	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 56.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3006	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.40% 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	0.44	0/870	0.68	0/1174
1	B	0.49	0/872	0.66	0/1177
1	C	0.46	0/929	0.66	0/1252
All	All	0.46	0/2671	0.67	0/3603

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 [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	853	0	878	18	0
1	B	843	0	884	42	0
1	C	888	0	951	42	0
2	A	133	0	0	0	0
2	B	133	0	0	3	0
2	C	156	0	0	1	0
All	All	3006	0	2713	64	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 12.

All (64) 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:B:421[B]:LYS:HG3	2:B:687:HOH:O	1.66	0.93
1:B:468[A]:GLN:HE21	1:B:469:LYS:H	1.21	0.89
1:C:388[B]:ASN:OD1	1:C:396:ASN:OD1	1.91	0.87
1:B:458:ILE:HD13	1:C:458[A]:ILE:HD11	1.59	0.85
1:B:488:LEU:HD11	1:C:491[B]:ILE:HD12	1.59	0.84
1:B:388:ASN:HB3	1:B:396:ASN:HD21	1.43	0.83
1:B:427:PHE:C	1:C:414[B]:LEU:HD23	1.98	0.83
1:B:418:LYS:HZ1	1:C:397:THR:HG21	1.47	0.80
1:B:418:LYS:NZ	1:C:397:THR:HG21	1.99	0.78
1:A:392:ASN:HD21	1:C:409[B]:SER:H	1.30	0.77
1:A:392:ASN:HD21	1:C:409[A]:SER:H	1.30	0.77
1:B:458:ILE:CD1	1:C:458[A]:ILE:HD11	2.14	0.77
1:B:418:LYS:HZ1	1:C:397:THR:CG2	1.98	0.76
1:B:418:LYS:NZ	1:C:397:THR:CG2	2.51	0.72
1:B:421[A]:LYS:HD3	1:B:422[A]:LYS:HE3	1.72	0.72
1:C:388[B]:ASN:HD21	1:C:396:ASN:HD21	1.39	0.69
1:C:385:GLY:N	1:C:401[B]:SER:HG	1.90	0.69
1:C:439[A]:LYS:HG3	2:C:585:HOH:O	1.92	0.68
1:B:458:ILE:HD13	1:C:458[A]:ILE:CD1	2.23	0.68
1:B:439:LYS:O	1:B:440[A]:LEU:HD23	1.97	0.63
1:B:439:LYS:C	1:B:440[A]:LEU:HD23	2.19	0.62
1:A:470[B]:ILE:HD13	1:C:463:ILE:HB	1.81	0.61
1:A:470[B]:ILE:CD1	1:C:463:ILE:HB	2.32	0.59
1:A:495:VAL:HG23	1:C:495:VAL:HG21	1.84	0.59
1:B:488:LEU:CD1	1:C:491[B]:ILE:HD12	2.30	0.59
1:B:458:ILE:CD1	1:C:458[A]:ILE:CD1	2.79	0.59
1:C:488:LEU:O	1:C:491[A]:ILE:HG22	2.04	0.56
1:C:388[B]:ASN:HD21	1:C:396:ASN:ND2	2.03	0.55
1:A:435:LEU:HB2	1:B:440[A]:LEU:CD2	2.37	0.54
1:B:395:ASN:ND2	2:B:652:HOH:O	2.40	0.54
1:C:388[B]:ASN:CG	1:C:396:ASN:OD1	2.46	0.54
1:B:427:PHE:O	1:C:414[B]:LEU:HD23	2.07	0.54
1:A:483[A]:VAL:HG23	1:B:483[A]:VAL:HG22	1.90	0.54
1:A:492:GLU:HB2	1:B:491:ILE:HD13	1.91	0.53
1:A:392:ASN:ND2	1:C:409[A]:SER:H	2.04	0.52
1:B:390:THR:HG23	1:B:396:ASN:HD22	1.74	0.52
1:A:392:ASN:ND2	1:C:409[B]:SER:H	2.03	0.52
1:A:436:LYS:HE3	1:C:431:ARG:HB2	1.92	0.52
1:B:466:GLY:O	1:B:468[B]:GLN:HG3	2.10	0.51
1:A:436:LYS:HE3	1:C:431:ARG:CB	2.40	0.51
1:B:427:PHE:HB3	1:C:414[B]:LEU:HD21	1.93	0.51
1:A:435:LEU:O	1:B:440[A]:LEU:HD22	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:LEU:HB2	1:B:440[A]:LEU:HD22	1.94	0.49
1:B:421[A]:LYS:HD2	2:B:687:HOH:O	2.12	0.49
1:B:427:PHE:HB3	1:C:414[B]:LEU:CD2	2.43	0.48
1:C:385:GLY:N	1:C:401[B]:SER:OG	2.46	0.48
1:B:438:ILE:HG22	1:B:440[A]:LEU:HD21	1.95	0.47
1:A:470[B]:ILE:CD1	1:C:463:ILE:HD12	2.45	0.46
1:B:428:ASP:N	1:C:414[B]:LEU:HD23	2.31	0.46
1:B:438:ILE:HG22	1:B:440[A]:LEU:CD2	2.46	0.46
1:B:428:ASP:CA	1:C:414[B]:LEU:HD22	2.45	0.46
1:B:438:ILE:CG2	1:B:440[A]:LEU:HD21	2.47	0.45
1:A:492:GLU:O	1:A:495:VAL:HG12	2.17	0.44
1:B:428:ASP:N	1:C:414[B]:LEU:CD2	2.81	0.44
1:B:468[A]:GLN:NE2	1:B:469:LYS:H	2.03	0.44
1:A:407:LEU:O	1:B:391:VAL:HA	2.19	0.42
1:B:418:LYS:HZ2	1:C:397:THR:CG2	2.30	0.42
1:B:423:ASN:ND2	1:C:397:THR:OG1	2.52	0.42
1:A:495:VAL:HG22	1:B:495:VAL:HG21	2.02	0.41
1:B:458:ILE:HD11	1:C:458[A]:ILE:HD11	1.97	0.41
1:C:388[B]:ASN:ND2	1:C:396:ASN:OD1	2.54	0.41
1:C:395:ASN:O	1:C:397:THR:HG23	2.21	0.41
1:B:418:LYS:NZ	1:C:397:THR:HG23	2.31	0.41

There are no symmetry-related clashes.

## 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	116/163 (71%)	113 (97%)	3 (3%)	0	100	100
1	B	117/163 (72%)	116 (99%)	1 (1%)	0	100	100
1	C	124/163 (76%)	122 (98%)	2 (2%)	0	100	100
All	All	357/489 (73%)	351 (98%)	6 (2%)	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	96/133 (72%)	96 (100%)	0	100	100
1	B	97/133 (73%)	95 (98%)	2 (2%)	59	16
1	C	104/133 (78%)	104 (100%)	0	100	100
All	All	297/399 (74%)	295 (99%)	2 (1%)	93	58

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	421[A]	LYS
1	B	421[B]	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	392	ASN
1	B	396	ASN
1	B	423	ASN
1	C	454	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

### 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	114/163 (69%)	1.06	13 (11%) 6 7	6, 10, 40, 54	0
1	B	111/163 (68%)	0.87	12 (10%) 6 8	6, 9, 26, 43	0
1	C	114/163 (69%)	1.06	12 (10%) 7 9	6, 9, 34, 52	0
All	All	339/489 (69%)	1.00	37 (10%) 6 8	6, 9, 36, 54	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	495	VAL	16.6
1	C	496	LYS	13.8
1	A	385	GLY	13.4
1	C	498	GLN	13.2
1	C	385	GLY	11.2
1	C	497	GLU	10.8
1	A	496	LYS	10.7
1	C	495	VAL	9.9
1	B	495	VAL	9.7
1	B	385	GLY	9.4
1	B	493	THR	9.2
1	B	491	ILE	7.6
1	A	497	GLU	7.0
1	A	386	GLY	6.9
1	C	494	GLU	6.5
1	A	498	GLN	6.2
1	B	492	GLU	5.9
1	A	491	ILE	5.9
1	C	491[A]	ILE	5.3
1	B	494	GLU	5.2
1	A	494	GLU	5.1
1	A	493	THR	4.8
1	A	492	GLU	4.4

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Mol	Chain	Res	Type	RSRZ
1	C	492	GLU	4.3
1	A	490	LYS	4.1
1	B	401	SER	4.0
1	B	386	GLY	3.4
1	A	421	LYS	3.1
1	C	493	THR	3.0
1	A	489	LYS	2.9
1	B	489	LYS	2.8
1	C	386	GLY	2.7
1	B	400	SER	2.4
1	C	414[A]	LEU	2.2
1	C	388[A]	ASN	2.1
1	B	490	LYS	2.1
1	B	454	ASN	2.1

## 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.