



# Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 2OSV  
Title : Crystal Structure of ZnuA from E. coli  
Authors : Li, H.; Jogl, G.  
Deposited on : 2007-02-06  
Resolution : 1.75 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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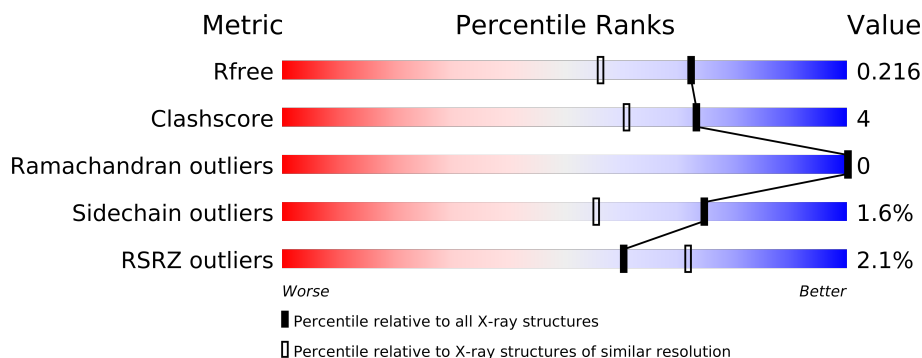
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.75 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))
$R_{free}$	66092	1134 (1.76-1.76)
Clashscore	79885	1304 (1.76-1.76)
Ramachandran outliers	78287	1288 (1.76-1.76)
Sidechain outliers	78261	1288 (1.76-1.76)
RSRZ outliers	66119	1135 (1.76-1.76)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	284	
1	B	284	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4704 atoms, of which 0 are hydrogen and 0 are deuterium.

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 High-affinity zinc uptake system protein znuA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	262	Total	C	N	O	S	0	4	0
			2042	1303	346	384	9			
1	B	263	Total	C	N	O	S	0	6	0
			2064	1317	349	389	9			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

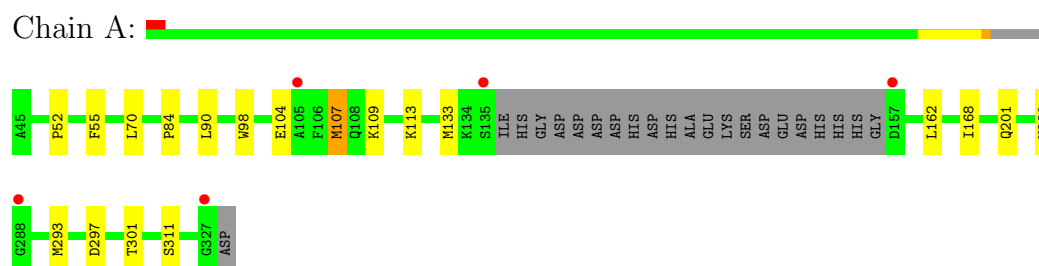
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	303	Total	O	0	0
			303	303		
3	B	293	Total	O	0	0
			293	293		

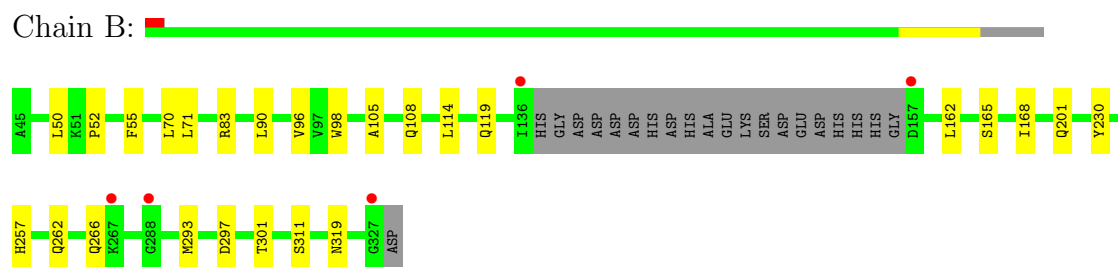
### 3 Residue-property plots

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 ( $RSRZ > 2$ ). 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: High-affinity zinc uptake system protein znuA



- Molecule 1: High-affinity zinc uptake system protein znuA



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	72.98Å 87.98Å 86.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.75 28.44 – 1.69	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-1.75) 99.5 (28.44-1.69)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.98 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.175 , 0.216 0.177 , 0.216	Depositor DCC
$R_{free}$ test set	3150 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	15.4	Xtriage
Anisotropy	0.325	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 37.4	EDS
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 62397 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4704	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.49 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.5373e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.57	0/2084	0.62	0/2825
1	B	0.56	0/2106	0.59	0/2855
All	All	0.57	0/4190	0.60	0/5680

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2042	0	2053	19	0
1	B	2064	0	2076	16	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	303	0	0	1	1
3	B	293	0	0	6	1
All	All	4704	0	4129	35	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (35) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:278:ARG:HE	1:A:278:ARG:HA	1.45	0.82
1:A:70:LEU:HD12	1:A:107[A]:MET:HE1	1.62	0.80
1:B:201:GLN:HG2	3:B:877:HOH:O	1.83	0.78
1:B:90[B]:LEU:HD13	1:B:96:VAL:HG21	1.73	0.70
1:A:287:ARG:NH1	3:A:748:HOH:O	2.23	0.69
1:A:278:ARG:NE	1:A:278:ARG:HA	2.15	0.62
1:B:70:LEU:HD11	1:B:90[A]:LEU:HD21	1.83	0.60
1:B:257:HIS:CD2	3:B:789:HOH:O	2.56	0.58
1:A:70:LEU:CD1	1:A:107[A]:MET:HE1	2.31	0.57
1:A:70:LEU:HD11	1:A:90:LEU:HD13	1.85	0.57
1:A:201:GLN:HG3	1:A:311:SER:OG	2.04	0.57
1:A:109:LYS:HG2	1:A:113:LYS:NZ	2.20	0.55
1:B:297:ASP:OD2	1:B:301[A]:THR:HG23	2.07	0.55
1:B:201:GLN:HG3	1:B:311:SER:OG	2.08	0.53
1:B:301[A]:THR:HG22	3:B:628:HOH:O	2.08	0.53
1:B:52:PRO:HA	1:B:55:PHE:CD2	2.44	0.52
1:A:109:LYS:HG2	1:A:113:LYS:HZ1	1.77	0.49
1:A:52:PRO:HA	1:A:55:PHE:CD2	2.48	0.49
1:B:90[B]:LEU:CD1	1:B:114:LEU:HD11	2.45	0.47
1:A:133:MET:HG2	1:A:230:TYR:CZ	2.49	0.47
1:B:105:ALA:HA	1:B:108:GLN:HG3	1.96	0.47
1:A:278:ARG:HE	1:A:278:ARG:CA	2.23	0.47
1:A:133:MET:HG2	1:A:230:TYR:CE1	2.50	0.46
1:B:165:SER:HB2	1:B:230:TYR:HB3	1.98	0.45
1:B:50:LEU:HD22	1:B:71:LEU:HD22	1.97	0.45
1:B:319:ASN:ND2	3:B:813:HOH:O	2.50	0.45
1:B:162:LEU:HB3	1:B:168:ILE:HG21	1.99	0.44
1:B:83:ARG:HD3	3:B:603:HOH:O	2.17	0.43
1:A:297:ASP:OD2	1:A:301:THR:HG23	2.19	0.43
1:A:84:PRO:HB2	3:B:830:HOH:O	2.18	0.43
1:A:107[A]:MET:HB3	1:A:107[A]:MET:HE3	1.63	0.43
1:A:107[A]:MET:HB2	1:A:107[A]:MET:HE2	1.63	0.42
1:A:162:LEU:HB3	1:A:168:ILE:HG21	2.02	0.42
1:A:104:GLU:O	1:A:107[A]:MET:HG2	2.20	0.41
1:B:262:GLN:HB2	1:B:266:GLN:HE21	1.86	0.41

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	Distance(Å)	Clash(Å)
3:A:730:HOH:O	3:B:842:HOH:O[2_654]	2.13	0.07

## 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	262/284 (92%)	259 (99%)	3 (1%)	0	100	100
1	B	265/284 (93%)	263 (99%)	2 (1%)	0	100	100
All	All	527/568 (93%)	522 (99%)	5 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

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	221/236 (94%)	216 (98%)	5 (2%)	63	36
1	B	224/236 (95%)	220 (98%)	4 (2%)	71	50
All	All	445/472 (94%)	436 (98%)	9 (2%)	75	43

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	TRP
1	A	107[A]	MET
1	A	107[B]	MET
1	A	278	ARG
1	A	293	MET
1	B	98	TRP
1	B	119	GLN
1	B	293[A]	MET
1	B	293[B]	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such



sidechains are listed below:

Mol	Chain	Res	Type
1	B	92	ASN
1	B	108	GLN
1	B	119	GLN
1	B	186	GLN
1	B	208	GLN
1	B	257	HIS
1	B	262	GLN
1	B	266	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	262/284 (92%)	-0.13	6 (2%) 57 68	9, 12, 20, 29	0
1	B	263/284 (92%)	-0.07	5 (1%) 64 74	7, 13, 23, 28	0
All	All	525/568 (92%)	-0.10	11 (2%) 60 72	7, 13, 22, 29	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	327	GLY	6.4
1	B	136	ILE	5.1
1	A	288	GLY	3.8
1	B	327	GLY	3.4
1	B	157	ASP	3.2
1	A	157	ASP	3.0
1	A	135	SER	2.8
1	A	105	ALA	2.4
1	B	288	GLY	2.2
1	B	267	LYS	2.1
1	A	287	ARG	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	B	602	1/1	0.03	-3.37	10,10,10,10	0
2	ZN	A	601	1/1	0.03	-4.28	12,12,12,12	0

## 6.5 Other polymers

There are no such residues in this entry.