



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 08:53 PM GMT

PDB ID : 1S1J  
Title : Crystal Structure of ZipA in complex with indoloquinolizin inhibitor 1  
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Deposited on : 2004-01-06  
Resolution : 2.18 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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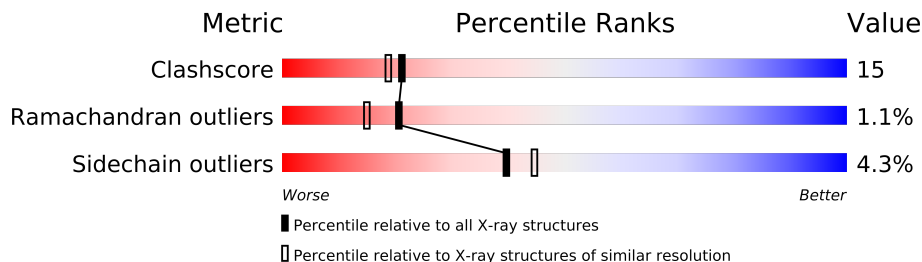
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : NOT EXECUTED  
EDS : NOT EXECUTED  
Percentile statistics : 21963  
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 2.18 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))
Clashscore	79885	4835 (2.20-2.16)
Ramachandran outliers	78287	4740 (2.20-2.16)
Sidechain outliers	78261	4741 (2.20-2.16)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	144	
1	B	144	

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 2384 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.

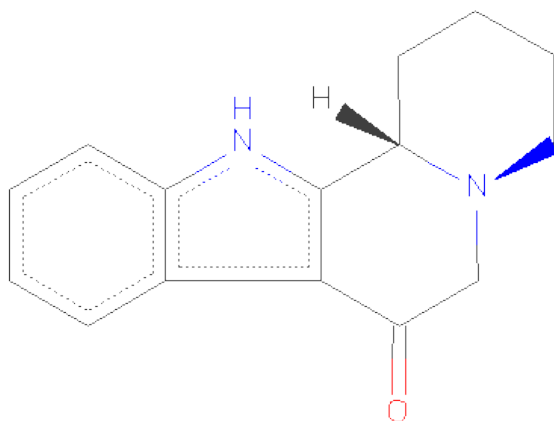
- Molecule 1 is a protein called Cell division protein zipA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	136	Total	C	N	O	S	0	0	0
			1063	673	181	201	8			
1	B	138	Total	C	N	O	S	0	0	0
			1077	680	184	205	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	97	LEU	ASN	CONFLICT	UNP P77173
B	97	LEU	ASN	CONFLICT	UNP P77173

- Molecule 2 is (12BS)-1,2,3,4,12,12B-HEXAHYDROINDOLO[2,3-A]QUINOLIZIN-7(6H)-ONE (three-letter code: IQZ) (formula: C<sub>15</sub>H<sub>16</sub>N<sub>2</sub>O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			18	15	2	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	87	Total 87	O 87	0	0
3	B	139	Total 139	O 139	0	0



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.79Å 39.46Å 70.51Å 90.00° 105.13° 90.00°	Depositor
Resolution (Å)	19.93 – 2.18	Depositor
% Data completeness (in resolution range)	98.1 (19.93-2.18)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.04	Depositor
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.192 , 0.236	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2384	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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

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	0.31	0/1084	0.55	0/1463
1	B	0.33	0/1098	0.59	0/1481
All	All	0.32	0/2182	0.57	0/2944

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	1063	0	1049	33	0
1	B	1077	0	1060	35	0
2	A	18	0	16	0	0
3	A	87	0	0	5	0
3	B	139	0	0	6	0
All	All	2384	0	2125	66	0

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 15.

All (66) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:93:ASP:O	1:B:97:LEU:HG	1.73	0.87
1:B:48:HIS:O	1:B:51:PRO:HD3	1.82	0.79
1:B:47:ARG:HH21	1:B:51:PRO:HG2	1.50	0.76
1:B:47:ARG:NH2	1:B:51:PRO:HG2	2.02	0.75
1:B:63:ASN:HD22	1:B:64:MET:H	1.35	0.75
1:B:71:ASP:HB3	1:B:74:MET:HB3	1.71	0.73
1:A:63:ASN:HD22	1:A:64:MET:N	1.86	0.73
1:B:63:ASN:HD22	1:B:64:MET:N	1.88	0.71
1:A:97:LEU:HD11	3:A:303:HOH:O	1.90	0.71
1:A:93:ASP:O	1:A:97:LEU:HG	1.92	0.70
1:B:8:GLU:HG3	1:B:87:GLN:NE2	2.09	0.68
1:B:137:ARG:HD3	1:B:137:ARG:O	1.96	0.66
1:B:29:LEU:HD11	3:B:239:HOH:O	1.96	0.65
1:B:47:ARG:HD3	3:B:204:HOH:O	1.95	0.65
1:A:63:ASN:HD22	1:A:64:MET:H	1.47	0.62
1:B:51:PRO:HA	3:B:228:HOH:O	2.00	0.61
1:B:46:HIS:HD2	1:B:60:SER:OG	1.84	0.61
1:A:76:ASP:O	1:A:77:PHE:HB3	2.00	0.61
1:B:50:SER:O	1:B:52:ASP:N	2.36	0.58
1:A:89:PRO:HD3	1:A:139:VAL:CG1	2.34	0.57
1:A:142:ALA:HA	3:A:353:HOH:O	2.04	0.57
1:B:28:LEU:O	1:B:32:ILE:HG13	2.06	0.55
1:A:37:PHE:HB3	1:A:45:TYR:HB3	1.88	0.55
1:A:50:SER:HB2	1:A:54:SER:O	2.08	0.54
1:A:38:ILE:HD12	3:A:335:HOH:O	2.08	0.53
1:B:30:ASN:HB3	3:B:225:HOH:O	2.09	0.52
1:B:107:HIS:HD2	3:B:234:HOH:O	1.93	0.52
1:A:58:LEU:HD13	1:A:97:LEU:HD22	1.91	0.52
1:A:71:ASP:OD1	1:A:73:GLU:HG2	2.09	0.52
1:B:74:MET:HG3	1:B:74:MET:O	2.10	0.51
1:A:48:HIS:HD2	1:A:55:GLY:O	1.94	0.50
1:A:46:HIS:HD2	1:A:60:SER:OG	1.94	0.50
1:A:107:HIS:CE1	1:A:111:GLU:OE2	2.65	0.50
1:B:37:PHE:HB3	1:B:45:TYR:HB3	1.94	0.49
1:B:49:LEU:HG	1:B:58:LEU:HD21	1.93	0.49
1:A:136:ILE:O	1:A:140:LYS:HG3	2.12	0.49
1:B:58:LEU:HD13	1:B:97:LEU:HD22	1.94	0.49
1:B:59:PHE:HB3	1:B:86:MET:HB2	1.94	0.48
1:A:97:LEU:HD23	3:A:311:HOH:O	2.13	0.48
1:A:63:ASN:ND2	1:A:65:VAL:H	2.11	0.47
1:A:66:LYS:HA	1:A:68:GLY:N	2.31	0.46
1:B:61:LEU:HD23	1:B:61:LEU:C	2.35	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:137:ARG:HD3	1:B:87:GLN:HG2	1.97	0.45
1:B:63:ASN:ND2	1:B:64:MET:N	2.62	0.45
1:A:41:ASP:O	1:A:46:HIS:HE1	2.00	0.45
1:B:91:TYR:O	1:B:97:LEU:HD11	2.16	0.44
1:A:87:GLN:HG2	3:A:363:HOH:O	2.17	0.44
1:A:52:ASP:OD1	1:A:54:SER:HB3	2.18	0.44
1:A:63:ASN:O	1:A:68:GLY:HA2	2.18	0.43
1:B:7:LYS:HG3	1:B:139:VAL:HG22	2.01	0.43
1:A:107:HIS:HE1	1:A:111:GLU:OE2	2.01	0.43
1:B:35:ALA:C	3:B:219:HOH:O	2.56	0.43
1:A:56:PRO:HB2	1:A:91:TYR:OH	2.20	0.42
1:B:7:LYS:HB2	1:B:8:GLU:H	1.66	0.42
1:B:88:VAL:HA	1:B:89:PRO:C	2.40	0.42
1:A:61:LEU:C	1:A:61:LEU:HD23	2.39	0.42
1:A:22:GLU:HB3	1:A:78:THR:HB	2.00	0.42
1:A:129:LEU:O	1:A:133:GLN:HG3	2.19	0.42
1:A:89:PRO:HD3	1:A:139:VAL:HG11	2.01	0.42
1:A:87:GLN:HE21	1:A:87:GLN:HB2	1.64	0.41
1:B:48:HIS:HD2	1:B:55:GLY:O	2.03	0.41
1:A:137:ARG:CD	1:B:87:GLN:HG2	2.51	0.40
1:A:90:SER:OG	1:A:91:TYR:N	2.53	0.40
1:B:46:HIS:CD2	1:B:60:SER:OG	2.70	0.40
1:B:11:ILE:O	1:B:85:PHE:HA	2.21	0.40
1:B:47:ARG:NE	1:B:51:PRO:HG3	2.37	0.40

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	134/144 (93%)	124 (92%)	8 (6%)	2 (2%)	15	10
1	B	136/144 (94%)	132 (97%)	3 (2%)	1 (1%)	30	26
All	All	270/288 (94%)	256 (95%)	11 (4%)	3 (1%)	21	15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	75	LYS
1	B	51	PRO
1	A	77	PHE

### 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	116/123 (94%)	110 (95%)	6 (5%)	32	35
1	B	117/123 (95%)	113 (97%)	4 (3%)	49	56
All	All	233/246 (95%)	223 (96%)	10 (4%)	40	45

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

Mol	Chain	Res	Type
1	A	38	ILE
1	A	41	ASP
1	A	52	ASP
1	A	63	ASN
1	A	73	GLU
1	A	87	GLN
1	B	51	PRO
1	B	63	ASN
1	B	74	MET
1	B	75	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	HIS
1	A	48	HIS
1	A	63	ASN
1	A	87	GLN
1	A	103	GLN
1	A	106	GLN

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Mol	Chain	Res	Type
1	A	107	HIS
1	A	127	GLN
1	B	18	HIS
1	B	33	GLN
1	B	34	GLN
1	B	46	HIS
1	B	48	HIS
1	B	63	ASN
1	B	87	GLN
1	B	107	HIS
1	B	143	ASN

### 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 ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	IQZ	A	301	-	21,21,21	1.79	4 (19%)	30,31,31	2.33	7 (23%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IQZ	A	301	-	-	0/0/23/23	0/0/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	IQZ	C12-C13	4.80	1.57	1.51
2	A	301	IQZ	C6-C1	2.50	1.42	1.36
2	A	301	IQZ	C5-C4	2.46	1.42	1.36
2	A	301	IQZ	C3-C2	2.25	1.45	1.41

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	IQZ	C9-C8-N7	9.33	110.98	106.87
2	A	301	IQZ	C8-C9-C13	4.31	123.88	122.51
2	A	301	IQZ	C9-C8-C10	-3.82	121.83	126.36
2	A	301	IQZ	C10-C8-N7	3.79	127.19	119.48
2	A	301	IQZ	C12-N11-C10	2.67	114.68	110.67
2	A	301	IQZ	C9-C3-C2	-2.58	106.10	107.82
2	A	301	IQZ	C16-C10-C8	-2.15	109.98	112.72

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.