



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 3W30  
Title : Structural basis for the recognition of Ubc13 by the *Shigella flexneri* effector OspI  
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Deposited on : 2012-12-07  
Resolution : 2.99 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](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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

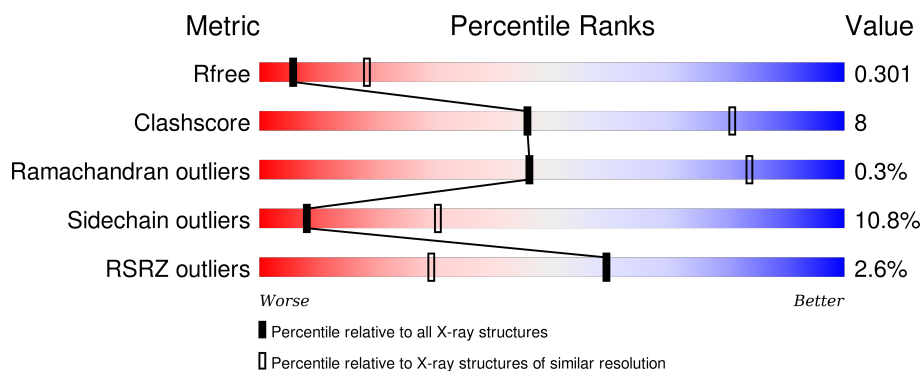
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.99 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	220	<div> <div>3%</div> <div>65%</div> <div>20%</div> <div>•</div> <div>13%</div> </div>
1	B	220	<div> <div>%</div> <div>66%</div> <div>20%</div> <div>•</div> <div>13%</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2980 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 ORF169b.

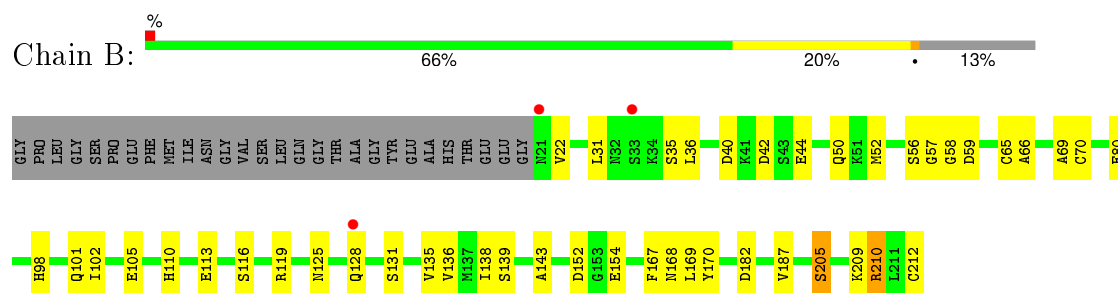
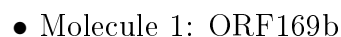
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	192	Total	C	N	O	S	0	0	0
			1490	927	255	298	10			
1	B	192	Total	C	N	O	S	0	0	0
			1490	927	255	298	10			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	EXPRESSION TAG	UNP Q8VSD5
A	-6	PRO	-	EXPRESSION TAG	UNP Q8VSD5
A	-5	LEU	-	EXPRESSION TAG	UNP Q8VSD5
A	-4	GLY	-	EXPRESSION TAG	UNP Q8VSD5
A	-3	SER	-	EXPRESSION TAG	UNP Q8VSD5
A	-2	PRO	-	EXPRESSION TAG	UNP Q8VSD5
A	-1	GLU	-	EXPRESSION TAG	UNP Q8VSD5
A	0	PHE	-	EXPRESSION TAG	UNP Q8VSD5
A	62	ALA	CYS	ENGINEERED MUTATION	UNP Q8VSD5
B	-7	GLY	-	EXPRESSION TAG	UNP Q8VSD5
B	-6	PRO	-	EXPRESSION TAG	UNP Q8VSD5
B	-5	LEU	-	EXPRESSION TAG	UNP Q8VSD5
B	-4	GLY	-	EXPRESSION TAG	UNP Q8VSD5
B	-3	SER	-	EXPRESSION TAG	UNP Q8VSD5
B	-2	PRO	-	EXPRESSION TAG	UNP Q8VSD5
B	-1	GLU	-	EXPRESSION TAG	UNP Q8VSD5
B	0	PHE	-	EXPRESSION TAG	UNP Q8VSD5
B	62	ALA	CYS	ENGINEERED MUTATION	UNP Q8VSD5



- Molecule 1: ORF169b



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.58Å 67.35Å 84.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.99 28.45 – 2.99	Depositor EDS
% Data completeness (in resolution range)	98.8 (30.00-2.99) 98.8 (28.45-2.99)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.91 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.230 , 0.303 0.227 , 0.301	Depositor DCC
$R_{free}$ test set	380 reflections (4.89%)	DCC
Wilson B-factor (Å <sup>2</sup> )	19.2	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 29.9	EDS
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.53$ , $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	3 of 8148 reflections (0.037%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2980	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	10.0	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 61.18 % 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 1.3609e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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 [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	0.72	0/1519	0.78	1/2049 (0.0%)
1	B	0.75	1/1519 (0.1%)	0.75	0/2049
All	All	0.74	1/3038 (0.0%)	0.77	1/4098 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	65	CYS	CB-SG	-6.21	1.71	1.82

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	ASP	CB-CG-OD1	5.38	123.14	118.30

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	1490	0	1428	22	0
1	B	1490	0	1428	23	0
All	All	2980	0	2856	45	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 8.

The worst 5 of 45 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:98:HIS:O	1:B:102:ILE:HD12	1.76	0.85
1:A:90:TYR:OH	1:A:199:HIS:ND1	2.11	0.83
1:B:119:ARG:HG2	1:B:119:ARG:HH11	1.48	0.77
1:B:42:ASP:OD2	1:B:119:ARG:NH2	2.14	0.76
1:B:31:LEU:HD11	1:B:44:GLU:HG2	1.69	0.75

There are no symmetry-related clashes.

## 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	190/220 (86%)	169 (89%)	21 (11%)	0	100	100
1	B	190/220 (86%)	175 (92%)	14 (7%)	1 (0%)	34	76
All	All	380/440 (86%)	344 (90%)	35 (9%)	1 (0%)	46	84

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	57	GLY

### 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	166/186 (89%)	144 (87%)	22 (13%)	5	21
1	B	166/186 (89%)	152 (92%)	14 (8%)	14	45
All	All	332/372 (89%)	296 (89%)	36 (11%)	8	30

5 of 36 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	176	SER
1	A	203	SER
1	B	187	VAL
1	A	196	GLU
1	A	210	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	ASN
1	B	168	ASN
1	B	32	ASN
1	A	125	ASN
1	B	98	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

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 [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<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	192/220 (87%)	0.01	7 (3%)	46 20	2, 10, 22, 28	0
1	B	192/220 (87%)	-0.02	3 (1%)	74 47	2, 9, 20, 24	0
All	All	384/440 (87%)	-0.01	10 (2%)	59 29	2, 10, 21, 28	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	21	ASN	2.9
1	A	129	ASN	2.6
1	B	33	SER	2.6
1	B	21	ASN	2.6
1	A	212	CYS	2.5

### 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.