



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 4ZV0  
Title : Structure of Tse6 in complex with Tsi6  
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Deposited on : 2015-05-18  
Resolution : 1.40 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

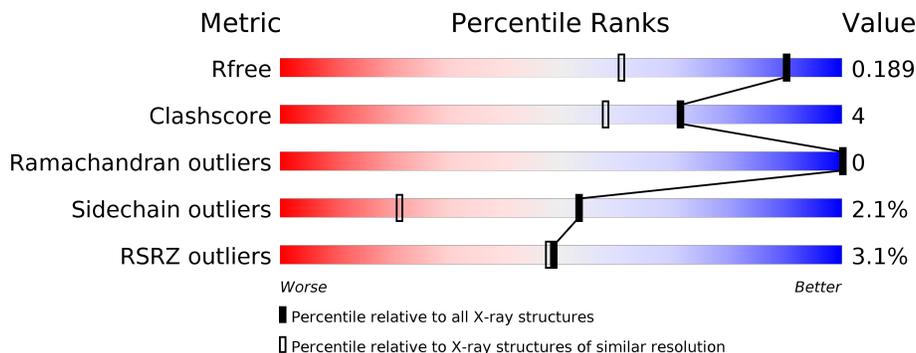
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.40 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))
$R_{free}$	130704	1714 (1.40-1.40)
Clashscore	141614	1812 (1.40-1.40)
Ramachandran outliers	138981	1763 (1.40-1.40)
Sidechain outliers	138945	1762 (1.40-1.40)
RSRZ outliers	127900	1674 (1.40-1.40)

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 respectively. 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	
2	B	94	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	IOD	B	103	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4040 atoms, of which 1820 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 antibacterial effector secreted protein (type VI secretion system).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	Se			
1	A	137	2112	680	1028	194	208	2	0	1	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	268	MSE	-	initiating methionine	UNP Q9I739
A	269	GLY	-	expression tag	UNP Q9I739
A	270	SER	-	expression tag	UNP Q9I739
A	271	SER	-	expression tag	UNP Q9I739
A	272	HIS	-	expression tag	UNP Q9I739
A	273	HIS	-	expression tag	UNP Q9I739
A	274	HIS	-	expression tag	UNP Q9I739
A	275	HIS	-	expression tag	UNP Q9I739
A	276	HIS	-	expression tag	UNP Q9I739
A	277	HIS	-	expression tag	UNP Q9I739
A	278	SER	-	expression tag	UNP Q9I739
A	279	GLN	-	expression tag	UNP Q9I739
A	280	ASP	-	expression tag	UNP Q9I739
A	281	PRO	-	expression tag	UNP Q9I739

- Molecule 2 is a protein called Tse6-binding/Tse6 immunity protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	Se			
2	B	94	1552	483	792	136	140	1	0	2	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MSE	-	initiating methionine	UNP Q9I740

- Molecule 3 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	5	Total I 5 5	0	0
3	A	5	Total I 5 5	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	225	Total O 225 225	0	0
4	B	141	Total O 141 141	0	0

### 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: antibacterial effector secreted protein (type VI secretion system)



- Molecule 2: Tse6-binding/Tse6 immunity protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.21Å 83.21Å 83.76Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.26 – 1.40 37.26 – 1.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (37.26-1.40) 100.0 (37.26-1.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.54 (at 1.40Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.146 , 0.164 0.170 , 0.189	Depositor DCC
$R_{free}$ test set	2022 reflections (3.13%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	12.9	Xtrriage
Anisotropy	0.280	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.42 , 50.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.038 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4040	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.89% 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](#)

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

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	0.91	3/1110 (0.3%)	1.09	5/1506 (0.3%)
2	B	0.84	2/784 (0.3%)	0.90	0/1060
All	All	0.88	5/1894 (0.3%)	1.02	5/2566 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	287	ARG	CG-CD	7.63	1.71	1.51
1	A	310	TYR	CE2-CZ	-5.34	1.31	1.38
2	B	9	ARG	CG-CD	-5.34	1.38	1.51
1	A	287	ARG	CZ-NH1	5.32	1.40	1.33
2	B	31	GLU	CD-OE2	-5.17	1.20	1.25

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	287	ARG	NE-CZ-NH1	23.03	131.81	120.30
1	A	287	ARG	NE-CZ-NH2	-13.59	113.50	120.30
1	A	287	ARG	CD-NE-CZ	7.26	133.77	123.60
1	A	303[A]	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	303[B]	ASP	CB-CG-OD1	5.20	122.98	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	1084	1028	1036	8	0
2	B	760	792	782	4	0
3	A	5	0	0	0	0
3	B	5	0	0	2	0
4	A	225	0	0	5	1
4	B	141	0	0	3	0
All	All	2220	1820	1818	13	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 4.

The worst 5 of 13 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)
3:B:103:IOD:I	4:B:320:HOH:O	2.45	1.02
1:A:393:PRO:O	4:A:601:HOH:O	1.87	0.92
4:A:655:HOH:O	3:B:103:IOD:I	2.60	0.90
2:B:44:ARG:NH1	4:B:201:HOH:O	2.14	0.80
1:A:385:GLU:CD	4:A:604:HOH:O	2.28	0.70

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)
4:A:744:HOH:O	4:A:764:HOH:O 4_665]	2.10	0.10

## 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	134/163 (82%)	133 (99%)	1 (1%)	0	100	100
2	B	94/94 (100%)	92 (98%)	2 (2%)	0	100	100
All	All	228/257 (89%)	225 (99%)	3 (1%)	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	115/134 (86%)	114 (99%)	1 (1%)	78	58
2	B	82/80 (102%)	79 (96%)	3 (4%)	34	7
All	All	197/214 (92%)	193 (98%)	4 (2%)	53	23

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	394	THR
2	B	32	LYS
2	B	44	ARG
2	B	65	ASP

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](#)

Of 10 ligands modelled in this entry, 10 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.

No monomer is involved in short contacts.

## 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	135/163 (82%)	-0.19	6 (4%) 34 34	9, 15, 33, 47	0
2	B	93/94 (98%)	-0.40	1 (1%) 80 79	7, 11, 25, 31	0
All	All	228/257 (88%)	-0.27	7 (3%) 49 48	7, 13, 29, 47	0

The worst 5 of 7 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	394	THR	7.0
1	A	393	PRO	4.8
1	A	396	ASP	3.3
1	A	395	ASP	3.0
1	A	397	PHE	2.7

### 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](#)

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. 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	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	IOD	A	501	1/1	0.79	0.26	10,10,10,10	1
3	IOD	B	102	1/1	0.82	0.14	56,56,56,56	1
3	IOD	A	503	1/1	0.93	0.08	29,29,29,29	1
3	IOD	B	104	1/1	0.97	0.05	25,25,25,25	1
3	IOD	A	504	1/1	0.97	0.06	18,18,18,18	1
3	IOD	A	502	1/1	0.97	0.05	29,29,29,29	1
3	IOD	A	505	1/1	0.97	0.06	25,25,25,25	1
3	IOD	B	101	1/1	0.99	0.04	15,15,15,15	1
3	IOD	B	105	1/1	0.99	0.04	22,22,22,22	1
3	IOD	B	103	1/1	0.99	0.05	28,28,28,28	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.