



# wwPDB X-ray Structure Validation Summary Report i

Feb 27, 2014 – 11:32 PM GMT

PDB ID : 3AQP  
Title : Crystal structure of SecDF, a translocon-associated membrane protein, from *Thermus thermophilus*  
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Deposited on : 2010-11-16  
Resolution : 3.30 Å (reported)

This is a wwPDB 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/ValidationPDFNotes.html>

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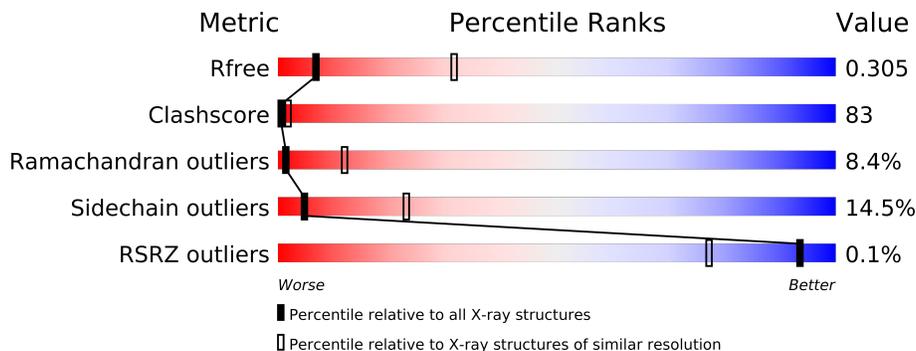
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.30 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	1341 (3.40-3.20)
Clashscore	79885	1696 (3.40-3.20)
Ramachandran outliers	78287	1664 (3.40-3.20)
Sidechain outliers	78261	1662 (3.40-3.20)
RSRZ outliers	66119	1342 (3.40-3.20)

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	741	
1	B	741	

## 2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 11028 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 Probable SecDF protein-export membrane protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	719	5514	3571	947	990	6	0	0	0
1	B	719	5514	3571	947	990	6	0	0	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ASP	ASN	ENGINEERED MUTATION	UNP Q5SKE6
A	736	HIS	-	EXPRESSION TAG	UNP Q5SKE6
A	737	HIS	-	EXPRESSION TAG	UNP Q5SKE6
A	738	HIS	-	EXPRESSION TAG	UNP Q5SKE6
A	739	HIS	-	EXPRESSION TAG	UNP Q5SKE6
A	740	HIS	-	EXPRESSION TAG	UNP Q5SKE6
A	741	HIS	-	EXPRESSION TAG	UNP Q5SKE6
B	2	ASP	ASN	ENGINEERED MUTATION	UNP Q5SKE6
B	736	HIS	-	EXPRESSION TAG	UNP Q5SKE6
B	737	HIS	-	EXPRESSION TAG	UNP Q5SKE6
B	738	HIS	-	EXPRESSION TAG	UNP Q5SKE6
B	739	HIS	-	EXPRESSION TAG	UNP Q5SKE6
B	740	HIS	-	EXPRESSION TAG	UNP Q5SKE6
B	741	HIS	-	EXPRESSION TAG	UNP Q5SKE6





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	140.64Å 140.64Å 160.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.88 – 3.30 46.88 – 3.30	Depositor EDS
% Data completeness (in resolution range)	89.0 (46.88-3.30) 89.0 (46.88-3.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.68 (at 3.32Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6_289)	Depositor
R, $R_{free}$	0.298 , 0.319 0.285 , 0.305	Depositor DCC
$R_{free}$ test set	2125 reflections (5.08%)	DCC
Wilson B-factor (Å <sup>2</sup> )	106.4	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 319.7	EDS
Estimated twinning fraction	0.501 for h,-k,-l 0.407 for h,-k,-l	Xtriage
Reported twinning fraction	0.501 for h,-k,-l	Depositor
L-test for twinning	$\langle  L  \rangle = 0.36$ , $\langle L^2 \rangle = 0.19$	Xtriage
Outliers	0 of 41834 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11028	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 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	0/5608	0.67	1/7622 (0.0%)
1	B	0.44	0/5608	0.67	1/7622 (0.0%)
All	All	0.44	0/11216	0.67	2/15244 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	165	LEU	CA-CB-CG	5.09	127.02	115.30
1	B	155	LEU	CA-CB-CG	5.05	126.91	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 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	5514	0	5787	939	0
1	B	5514	0	5787	944	0
All	All	11028	0	11574	1880	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 83.

The worst 5 of 1880 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:41:GLY:O	1:A:94:GLY:HA2	1.32	1.27
1:B:41:GLY:O	1:B:94:GLY:HA2	1.31	1.24
1:B:683:ILE:HD11	1:B:702:PHE:HB2	1.23	1.20
1:B:541:GLU:HA	1:B:544:PHE:CD1	1.79	1.17
1:A:541:GLU:HA	1:A:544:PHE:CD1	1.80	1.16

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	715/741 (96%)	463 (65%)	193 (27%)	59 (8%)	1	12
1	B	715/741 (96%)	456 (64%)	198 (28%)	61 (8%)	1	11
All	All	1430/1482 (96%)	919 (64%)	391 (27%)	120 (8%)	1	12

5 of 120 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	PRO
1	A	76	ALA
1	A	331	LEU
1	A	410	SER
1	A	574	VAL

### 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	581/597 (97%)	496 (85%)	85 (15%)	5	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	581/597 (97%)	497 (86%)	84 (14%)	5	23
All	All	1162/1194 (97%)	993 (86%)	169 (14%)	5	23

5 of 169 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	680	LEU
1	B	80	ILE
1	B	670	SER
1	A	684	LEU
1	B	16	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	667	GLN
1	B	68	ASN
1	B	420	HIS
1	A	662	ASN
1	B	569	ASN

### 5.3.3 RNA [i](#)

There are no RNA chains 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

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	719/741 (97%)	-0.31	1 (0%) 93 77	31, 111, 181, 242	0
1	B	719/741 (97%)	-0.29	1 (0%) 93 77	33, 111, 184, 240	0
All	All	1438/1482 (97%)	-0.30	2 (0%) 93 77	31, 111, 183, 242	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	34	ARG	2.3
1	B	213	TYR	2.3

### 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

There are no ligands in this entry.

### 6.5 Other polymers

There are no such residues in this entry.