



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 13, 2024 – 09:53 AM EST

PDB ID : 4ODL  
Title : Structure of SlyD from *Thermus thermophilus* in complex with S2 peptide  
Authors : Quistgaard, E.M.; Low, C.; Nordlund, P.  
Deposited on : 2014-01-10  
Resolution : 2.92 Å(reported)

This is a Full wwPDB X-ray Structure 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

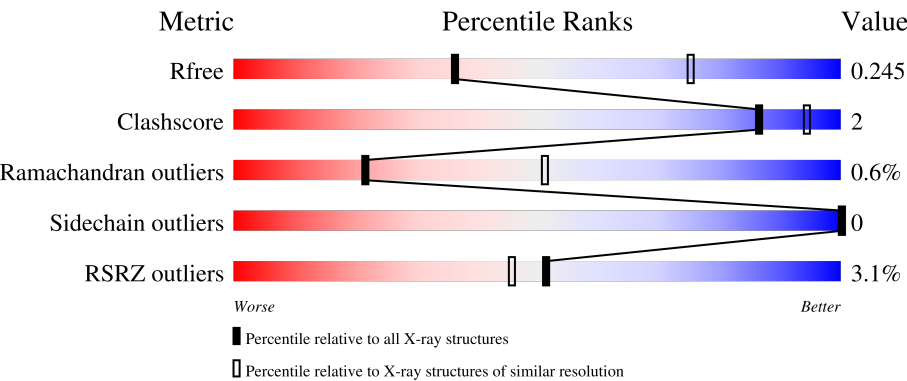
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:  
*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.92 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 <sub>free</sub>	164625	2797 (2.94-2.90)
Clashscore	180529	3049 (2.94-2.90)
Ramachandran outliers	177936	2981 (2.94-2.90)
Sidechain outliers	177891	2983 (2.94-2.90)
RSRZ outliers	164620	2799 (2.94-2.90)

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 of 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	158	<div><div>%</div><div><div></div><div>91%</div><div>• 5%</div></div></div>
1	B	158	<div><div>%</div><div><div></div><div>91%</div><div>• 5%</div></div></div>
2	C	16	<div><div>12%</div><div><div></div><div>44%</div><div>19%</div><div>38%</div></div></div>
2	D	16	<div><div>12%</div><div><div></div><div>81%</div><div>19%</div></div></div>
2	E	16	<div><div>12%</div><div><div></div><div>94%</div><div>6%</div></div></div>

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Mol	Chain	Length	Quality of chain
2	F	16	<div><div></div><div>6%</div><div>94%</div><div>6%</div></div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2791 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 Peptidyl-prolyl cis-trans isomerase SlyD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	150	Total	C	N	O	S	0	0	0
			1157	730	194	230	3			
1	B	150	Total	C	N	O	S	0	0	0
			1157	730	194	230	3			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	150	PRO	-	expression tag	UNP Q5SLE7
A	151	SER	-	expression tag	UNP Q5SLE7
A	152	GLY	-	expression tag	UNP Q5SLE7
A	153	HIS	-	expression tag	UNP Q5SLE7
A	154	HIS	-	expression tag	UNP Q5SLE7
A	155	HIS	-	expression tag	UNP Q5SLE7
A	156	HIS	-	expression tag	UNP Q5SLE7
A	157	HIS	-	expression tag	UNP Q5SLE7
A	158	HIS	-	expression tag	UNP Q5SLE7
B	150	PRO	-	expression tag	UNP Q5SLE7
B	151	SER	-	expression tag	UNP Q5SLE7
B	152	GLY	-	expression tag	UNP Q5SLE7
B	153	HIS	-	expression tag	UNP Q5SLE7
B	154	HIS	-	expression tag	UNP Q5SLE7
B	155	HIS	-	expression tag	UNP Q5SLE7
B	156	HIS	-	expression tag	UNP Q5SLE7
B	157	HIS	-	expression tag	UNP Q5SLE7
B	158	HIS	-	expression tag	UNP Q5SLE7

- Molecule 2 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	10	Total	C	N	O	S	0	0	1
			73	51	12	9	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	16	Total	C	N	O	S	0	0	1
			132	90	23	18	1			
2	E	16	Total	C	N	O	S	0	0	1
			132	90	23	18	1			
2	F	16	Total	C	N	O	S	0	0	1
			132	90	23	18	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	35	NH2	-	amidation	UNP P0A7V0
D	35	NH2	-	amidation	UNP P0A7V0
E	35	NH2	-	amidation	UNP P0A7V0
F	35	NH2	-	amidation	UNP P0A7V0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

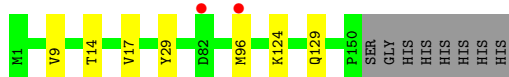
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is water.

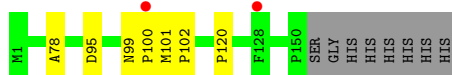
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		
4	B	4	Total	O	0	0
			4	4		

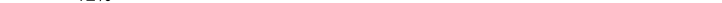


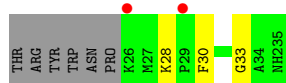
- Molecule 1: Peptidyl-prolyl cis-trans isomerase SlyD



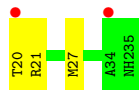
- Chain B:  91% 5%



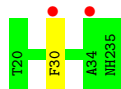
- Chain C: 



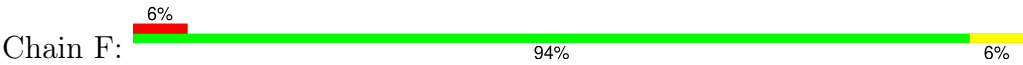
- Chain D:  12% 81% 19%



- Chain E:  12% 94% 6%



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.88Å 110.88Å 182.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.01 – 2.92 48.01 – 2.92	Depositor EDS
% Data completeness (in resolution range)	99.1 (48.01-2.92) 99.1 (48.01-2.92)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.83 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, $R_{free}$	0.210 , 0.224 0.230 , 0.245	Depositor DCC
$R_{free}$ test set	756 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	84.8	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 48.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	2791	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	74.0	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 21.33 % 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 7.1813e-03.*

<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: NH2, CL

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	0.22	0/1185	0.41	0/1615
1	B	0.21	0/1185	0.41	0/1615
2	C	0.26	0/74	0.60	0/96
2	D	0.31	0/137	0.56	0/184
2	E	0.23	0/137	0.38	0/184
2	F	0.26	0/137	0.51	0/184
All	All	0.22	0/2855	0.43	0/3878

There are no bond length outliers.

There are no bond angle outliers.

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	1157	0	1104	4	0
1	B	1157	0	1104	6	0
2	C	73	0	78	2	0
2	D	132	0	130	2	0
2	E	132	0	130	1	0
2	F	132	0	130	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
4	A	2	0	0	0	0
4	B	4	0	0	0	0
All	All	2791	0	2676	11	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 2.

All (11) 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:A:14:THR:HB	1:A:129:GLN:HB3	1.88	0.56
1:A:96:MET:HG2	2:D:27:MET:HB3	1.87	0.55
1:A:17:VAL:HG11	1:A:124:LYS:HD3	1.92	0.51
1:B:120:PRO:HG2	2:E:30:PHE:HE2	1.78	0.48
1:B:95:ASP:OD1	1:B:99:ASN:N	2.49	0.46
2:D:20:THR:HA	2:D:21:ARG:HA	1.68	0.45
1:B:78:ALA:O	2:C:33:GLY:HA3	2.19	0.43
1:B:99:ASN:HA	1:B:100:PRO:HD3	1.81	0.42
1:B:95:ASP:HA	2:C:30:PHE:HB3	2.03	0.41
1:B:101:MET:HA	1:B:102:PRO:HD2	1.97	0.41
1:A:9:VAL:HB	1:A:29:TYR:CE2	2.56	0.40

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	148/158 (94%)	143 (97%)	5 (3%)	0	100	100
1	B	148/158 (94%)	146 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	8/16 (50%)	4 (50%)	3 (38%)	1 (12%)	0	0
2	D	14/16 (88%)	9 (64%)	5 (36%)	0	100	100
2	E	14/16 (88%)	12 (86%)	2 (14%)	0	100	100
2	F	14/16 (88%)	13 (93%)	0	1 (7%)	1	2
All	All	346/380 (91%)	327 (94%)	17 (5%)	2 (1%)	22	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	34	ALA
2	C	28	LYS

### 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	123/130 (95%)	123 (100%)	0	100	100
1	B	123/130 (95%)	123 (100%)	0	100	100
2	C	7/13 (54%)	7 (100%)	0	100	100
2	D	13/13 (100%)	13 (100%)	0	100	100
2	E	13/13 (100%)	13 (100%)	0	100	100
2	F	13/13 (100%)	13 (100%)	0	100	100
All	All	292/312 (94%)	292 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

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.

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 ⓘ

### 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	150/158 (94%)	-0.46	2 (1%) 74 69	30, 63, 103, 126	0
1	B	150/158 (94%)	-0.29	2 (1%) 74 69	28, 73, 122, 145	0
2	C	9/16 (56%)	1.51	2 (22%) 3 3	90, 117, 133, 144	0
2	D	15/16 (93%)	1.12	2 (13%) 8 8	52, 98, 137, 139	0
2	E	15/16 (93%)	0.21	2 (13%) 8 8	54, 63, 98, 120	0
2	F	15/16 (93%)	0.28	1 (6%) 25 22	50, 66, 109, 112	0
All	All	354/380 (93%)	-0.21	11 (3%) 51 46	28, 70, 122, 145	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	20	THR	5.5
2	C	29	PRO	4.3
2	E	30	PHE	3.2
2	D	34	ALA	3.1
2	C	26	LYS	2.8
2	F	34	ALA	2.8
2	E	34	ALA	2.8
1	A	96	MET	2.6
1	B	128	PHE	2.2
1	B	100	PRO	2.2
1	A	82	ASP	2.1

### 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 [i](#)

There are no monosaccharides 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	CL	A	201	1/1	0.98	0.04	50,50,50,50	0
3	CL	B	201	1/1	0.99	0.02	51,51,51,51	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.