



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 11:32 pm GMT

PDB ID : 4YTO  
Title : Structure of coiled-coil domain of SYCP1  
Authors : park, H.H.  
Deposited on : 2015-03-18  
Resolution : 2.00 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

<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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

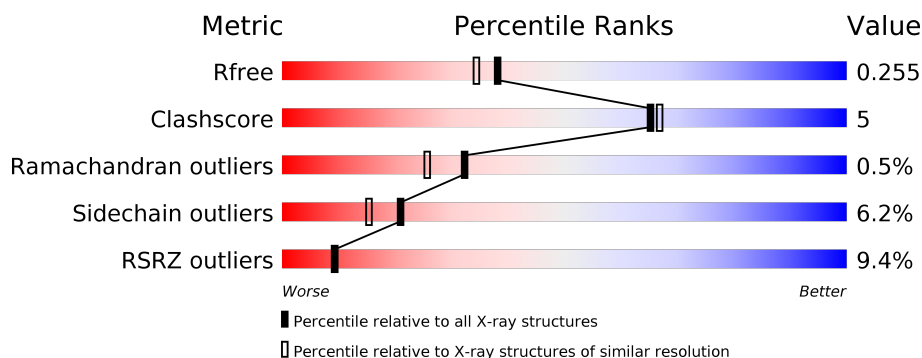
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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	148	<div> <div>7%</div> <div>51%</div> <div>9%</div> <div>•</div> <div>36%</div> </div>
1	B	148	<div> <div>5%</div> <div>55%</div> <div>9%</div> <div>•</div> <div>34%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1728 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.

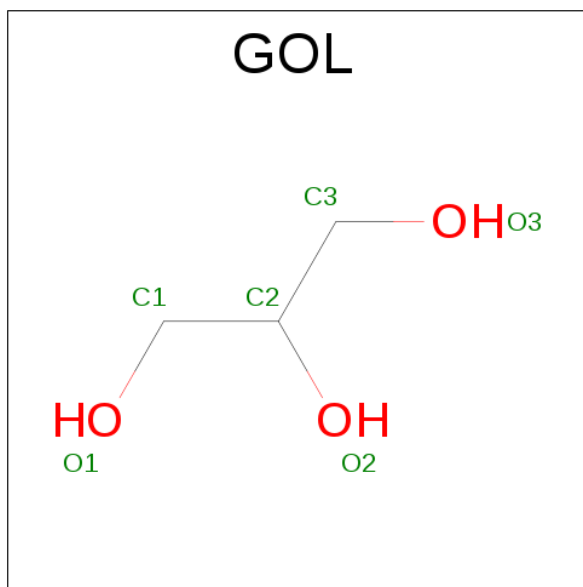
- Molecule 1 is a protein called Synaptonemal complex protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	95	Total	C	N	O	S	0	0	0
			772	482	132	155	3			
1	B	97	Total	C	N	O	S	0	0	0
			789	493	134	159	3			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	802	LEU	-	expression tag	UNP Q15431
A	803	GLU	-	expression tag	UNP Q15431
A	804	HIS	-	expression tag	UNP Q15431
A	805	HIS	-	expression tag	UNP Q15431
A	806	HIS	-	expression tag	UNP Q15431
A	807	HIS	-	expression tag	UNP Q15431
A	808	HIS	-	expression tag	UNP Q15431
A	809	HIS	-	expression tag	UNP Q15431
B	802	LEU	-	expression tag	UNP Q15431
B	803	GLU	-	expression tag	UNP Q15431
B	804	HIS	-	expression tag	UNP Q15431
B	805	HIS	-	expression tag	UNP Q15431
B	806	HIS	-	expression tag	UNP Q15431
B	807	HIS	-	expression tag	UNP Q15431
B	808	HIS	-	expression tag	UNP Q15431
B	809	HIS	-	expression tag	UNP Q15431

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	64	Total	O	0	0
			64	64		
3	B	91	Total	O	0	0
			91	91		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	119.61Å 119.61Å 83.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.19 – 2.00 28.19 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.5 (28.19-2.00) 97.6 (28.19-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.98 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.231 , 0.256 0.236 , 0.255	Depositor DCC
$R_{free}$ test set	2000 reflections (5.16%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.5	Xtriage
Anisotropy	0.320	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 49.7	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	0.93	EDS
Total number of atoms	1728	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.91% 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: GOL

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	1.54	9/776 (1.2%)	1.21	3/1032 (0.3%)
1	B	1.61	10/793 (1.3%)	1.12	4/1055 (0.4%)
All	All	1.58	19/1569 (1.2%)	1.17	7/2087 (0.3%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	768	GLU	CD-OE1	10.42	1.37	1.25
1	B	768	GLU	CD-OE2	7.81	1.34	1.25
1	A	754	SER	CB-OG	-7.80	1.32	1.42
1	B	754	SER	CB-OG	-7.76	1.32	1.42
1	B	698	GLU	CG-CD	7.30	1.62	1.51
1	B	743	SER	CB-OG	-7.10	1.33	1.42
1	B	698	GLU	CD-OE1	6.92	1.33	1.25
1	B	691	GLU	CB-CG	-6.91	1.39	1.52
1	B	691	GLU	CD-OE2	-6.71	1.18	1.25
1	A	691	GLU	CD-OE2	-6.71	1.18	1.25
1	A	739	GLU	CD-OE1	6.48	1.32	1.25
1	A	750	GLU	CD-OE2	6.46	1.32	1.25
1	B	735	TYR	CE1-CZ	-6.41	1.30	1.38
1	A	742	GLN	CB-CG	-5.99	1.36	1.52
1	A	715	GLU	CD-OE1	5.68	1.31	1.25
1	B	743	SER	CA-CB	-5.47	1.44	1.52
1	A	739	GLU	CB-CG	5.25	1.62	1.52
1	A	740	GLN	CG-CD	5.13	1.62	1.51
1	A	691	GLU	CG-CD	5.02	1.59	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	679	LEU	CA-CB-CG	6.37	129.96	115.30
1	B	746	ARG	NE-CZ-NH2	-6.33	117.14	120.30
1	A	739	GLU	OE1-CD-OE2	-6.03	116.06	123.30
1	A	746	ARG	NE-CZ-NH1	-5.41	117.60	120.30
1	A	716	LYS	CD-CE-NZ	5.18	123.63	111.70
1	B	698	GLU	CA-CB-CG	5.15	124.74	113.40
1	B	752	GLU	OE1-CD-OE2	-5.12	117.15	123.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-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 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	772	0	803	12	0
1	B	789	0	820	4	0
2	A	12	0	15	0	0
3	A	64	0	0	5	0
3	B	91	0	0	2	0
All	All	1728	0	1638	15	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 5.

All (15) 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:673:ILE:HG22	1:A:674:SER:H	1.43	0.82
1:A:767:LEU:HD12	3:A:1064:HOH:O	1.81	0.79
1:A:719:HIS:HB3	3:A:1054:HOH:O	1.87	0.73
1:A:673:ILE:HG22	1:A:674:SER:N	2.04	0.71
1:A:673:ILE:CG2	1:A:674:SER:H	2.11	0.63
1:A:740:GLN:CG	3:A:1052:HOH:O	2.59	0.50
1:A:740:GLN:HG2	3:A:1052:HOH:O	2.13	0.48
1:A:767:LEU:CD1	3:A:1064:HOH:O	2.50	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:706:LYS:NZ	3:B:905:HOH:O	2.49	0.44
1:A:715:GLU:HG3	1:B:725:ILE:HD13	2.00	0.43
1:B:687:VAL:O	1:B:691:GLU:HG2	2.19	0.43
1:A:701:LYS:HA	1:A:701:LYS:HD3	1.83	0.42
1:B:686:LYS:NZ	3:B:903:HOH:O	2.43	0.41
1:A:728:ARG:NH2	1:A:732:LEU:HD21	2.36	0.41
1:A:765:LYS:HB2	1:A:765:LYS:HE2	1.81	0.41

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	93/148 (63%)	90 (97%)	2 (2%)	1 (1%)	17	9
1	B	95/148 (64%)	95 (100%)	0	0	100	100
All	All	188/296 (64%)	185 (98%)	2 (1%)	1 (0%)	32	26

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	674	SER

### 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	87/138 (63%)	81 (93%)	6 (7%)	18	12
1	B	89/138 (64%)	84 (94%)	5 (6%)	25	19
All	All	176/276 (64%)	165 (94%)	11 (6%)	21	15

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	678	LEU
1	A	732	LEU
1	A	754	SER
1	A	761	LEU
1	A	762	SER
1	A	765	LYS
1	B	679	LEU
1	B	706	LYS
1	B	718	LYS
1	B	730	SER
1	B	769	ILE

Some 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 ⓘ

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 ⓘ

2 ligands are 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	GOL	A	901	-	5,5,5	0.37	0	5,5,5	1.07	0
2	GOL	A	902	-	5,5,5	1.18	0	5,5,5	1.36	1 (20%)

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	GOL	A	901	-	-	0/4/4/4	0/0/0/0
2	GOL	A	902	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	902	GOL	O1-C1-C2	-2.33	98.34	110.07

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	95/148 (64%)	0.45	11 (11%) 5 5	21, 39, 91, 108	0
1	B	97/148 (65%)	0.30	7 (7%) 16 16	20, 41, 81, 103	0
All	All	192/296 (64%)	0.38	18 (9%) 9 9	20, 40, 84, 108	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	768	GLU	7.2
1	B	767	LEU	5.8
1	A	767	LEU	5.2
1	A	764	LYS	4.8
1	A	673	ILE	4.8
1	A	766	GLN	4.7
1	B	769	ILE	4.6
1	B	766	GLN	4.4
1	B	673	ILE	3.7
1	A	765	LYS	3.6
1	A	726	GLU	3.5
1	A	732	LEU	3.4
1	B	765	LYS	3.2
1	A	729	ASP	3.0
1	A	731	GLU	2.9
1	A	762	SER	2.6
1	B	764	LYS	2.6
1	A	675	GLU	2.1

### 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GOL	A	902	6/6	0.86	0.16	-0.12	43,50,58,75	0
2	GOL	A	901	6/6	0.98	0.09	-0.38	31,33,36,41	0

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

There are no such residues in this entry.