



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 24, 2024 – 02:37 pm BST

PDB ID : 8PHX  
Title : Receiver Domain of the Hybrid Histidine Kinase Sln1 of *Candida albicans*  
Authors : Paredes-Martinez, F.; Casino, P.  
Deposited on : 2023-06-20  
Resolution : 1.50 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.

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) : 1.13  
EDS : 2.37.1  
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) : 2.37.1

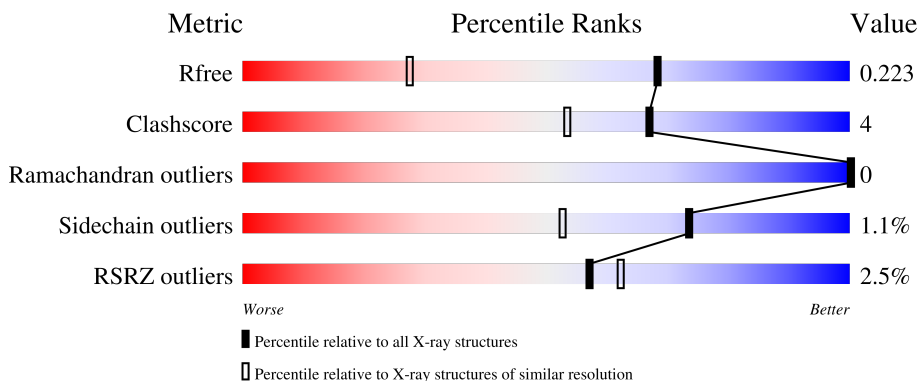
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.50 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	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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	132	 89% 7% . .
1	B	132	 88% 10% .
1	C	132	 92% 7% .
1	D	132	 92% 6% .

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4494 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 Histidine protein kinase SLN1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	128	996	633	165	190	8	0	4	0
1	B	129	1001	634	168	191	8	0	2	0
1	C	130	1011	640	167	196	8	0	4	0
1	D	129	1002	634	168	192	8	0	1	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1237	GLY	-	expression tag	UNP Q5A872
A	1238	PRO	-	expression tag	UNP Q5A872
A	1239	GLY	-	expression tag	UNP Q5A872
A	1253	LEU	SER	variant	UNP Q5A872
B	1237	GLY	-	expression tag	UNP Q5A872
B	1238	PRO	-	expression tag	UNP Q5A872
B	1239	GLY	-	expression tag	UNP Q5A872
B	1253	LEU	SER	variant	UNP Q5A872
C	1237	GLY	-	expression tag	UNP Q5A872
C	1238	PRO	-	expression tag	UNP Q5A872
C	1239	GLY	-	expression tag	UNP Q5A872
C	1253	LEU	SER	variant	UNP Q5A872
D	1237	GLY	-	expression tag	UNP Q5A872
D	1238	PRO	-	expression tag	UNP Q5A872
D	1239	GLY	-	expression tag	UNP Q5A872
D	1253	LEU	SER	variant	UNP Q5A872

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0

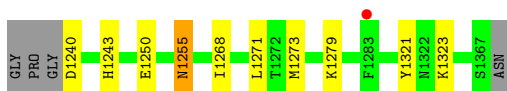
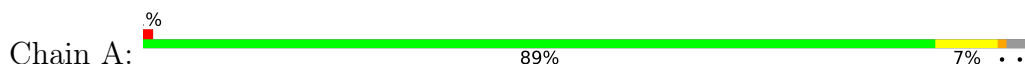
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	110	Total O 110 110	0	0
3	B	113	Total O 113 113	0	0
3	C	122	Total O 122 122	0	0
3	D	135	Total O 135 135	0	0

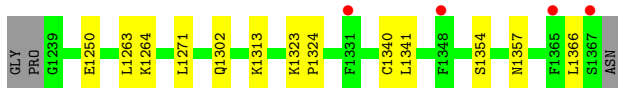
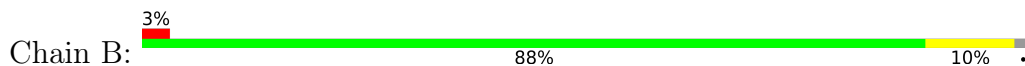
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Histidine protein kinase SLN1



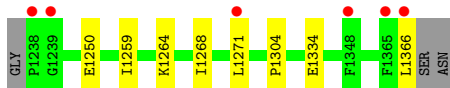
- Molecule 1: Histidine protein kinase SLN1



- Molecule 1: Histidine protein kinase SLN1



- Molecule 1: Histidine protein kinase SLN1



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.16Å 96.27Å 97.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.20 – 1.50 54.14 – 1.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (54.20-1.50) 100.0 (54.14-1.50)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.45 (at 1.50Å)	Xtrriage
Refinement program	REFMAC 5.8.0411	Depositor
R, $R_{free}$	0.193 , 0.215 0.203 , 0.223	Depositor DCC
$R_{free}$ test set	4801 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	17.8	Xtrriage
Anisotropy	0.129	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 39.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.000 for -h,l,k	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4494	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 81.91 % 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 3.1707e-07. 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, 0.333 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: MG

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	0.48	2/1018 (0.2%)	0.72	2/1373 (0.1%)
1	B	0.41	0/1017	0.69	0/1368
1	C	0.38	0/1033	0.67	1/1392 (0.1%)
1	D	0.38	0/1015	0.72	0/1365
All	All	0.41	2/4083 (0.0%)	0.70	3/5498 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1255[A]	ASN	C-O	6.41	1.35	1.23
1	A	1255[B]	ASN	C-O	6.41	1.35	1.23

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1238	PRO	N-CA-CB	5.96	110.45	103.30
1	A	1255[A]	ASN	CA-C-O	5.07	130.74	120.10
1	A	1255[B]	ASN	CA-C-O	5.07	130.74	120.10

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	996	0	1026	11	0
1	B	1001	0	1034	11	0
1	C	1011	0	1037	4	1
1	D	1002	0	1033	8	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	110	0	0	4	0
3	B	113	0	0	5	0
3	C	122	0	0	1	1
3	D	135	0	0	4	2
All	All	4494	0	4130	32	2

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 32 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:1313:LYS:HD2	3:B:1540:HOH:O	1.35	1.27
1:A:1240:ASP:HA	3:A:1501:HOH:O	1.43	1.18
1:D:1259:ILE:HG13	3:D:1561:HOH:O	1.77	0.83
1:D:1268:ILE:HG21	1:D:1271:LEU:HD21	1.69	0.73
1:D:1259:ILE:CG1	3:D:1561:HOH:O	2.35	0.72

All (2) 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)
1:C:1367:SER:C	3:D:1501:HOH:O[1_455]	1.62	0.58
3:C:1508:HOH:O	3:D:1626:HOH:O[1_455]	2.14	0.06

## 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	130/132 (98%)	127 (98%)	3 (2%)	0	100	100
1	B	129/132 (98%)	126 (98%)	3 (2%)	0	100	100
1	C	132/132 (100%)	129 (98%)	3 (2%)	0	100	100
1	D	128/132 (97%)	123 (96%)	5 (4%)	0	100	100
All	All	519/528 (98%)	505 (97%)	14 (3%)	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	114/117 (97%)	113 (99%)	1 (1%)	78	61
1	B	114/117 (97%)	112 (98%)	2 (2%)	59	30
1	C	116/117 (99%)	115 (99%)	1 (1%)	78	61
1	D	114/117 (97%)	113 (99%)	1 (1%)	78	61
All	All	458/468 (98%)	453 (99%)	5 (1%)	73	53

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1250	GLU
1	B	1250	GLU
1	B	1366	LEU
1	C	1250	GLU
1	D	1250	GLU

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

Mol	Chain	Res	Type
1	B	1357	ASN
1	C	1255	ASN
1	D	1320	GLN
1	D	1255	ASN
1	B	1255	ASN

### 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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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	128/132 (96%)	0.05	1 (0%) 86 89	12, 20, 34, 41	0
1	B	129/132 (97%)	0.07	4 (3%) 49 54	12, 19, 36, 43	0
1	C	130/132 (98%)	-0.08	2 (1%) 73 78	13, 18, 29, 34	0
1	D	129/132 (97%)	0.08	6 (4%) 31 34	13, 19, 32, 38	0
All	All	516/528 (97%)	0.03	13 (2%) 57 62	12, 19, 32, 43	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1367	SER	3.5
1	B	1365	PHE	3.3
1	C	1238	PRO	3.2
1	D	1239	GLY	2.8
1	B	1331	PHE	2.8

### 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 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
2	MG	D	1401	1/1	0.98	0.04	15,15,15,15	0
2	MG	B	1401	1/1	0.99	0.05	14,14,14,14	0
2	MG	C	1401	1/1	0.99	0.05	15,15,15,15	0
2	MG	A	1401	1/1	0.99	0.05	14,14,14,14	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.