



# wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 1H3Z  
Title : SOLUTION STRUCTURE OF A PWWP DOMAIN FROM SCHIZOSACCHAROMYCES POMBE  
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Deposited on : 2002-09-23

This is a wwPDB NMR 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  
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>  
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:

Cyrange : Kirchner and Güntert (2011)  
NmrClust : Kelley et al. (1996)  
MolProbity : 4.02b-467  
Mogul : unknown  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
RCI : v\_1n\_11\_5\_13\_A (Berjanski et al., 2005)  
PANAV : Wang et al. (2010)  
ShiftChecker : rb-20027457  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027457

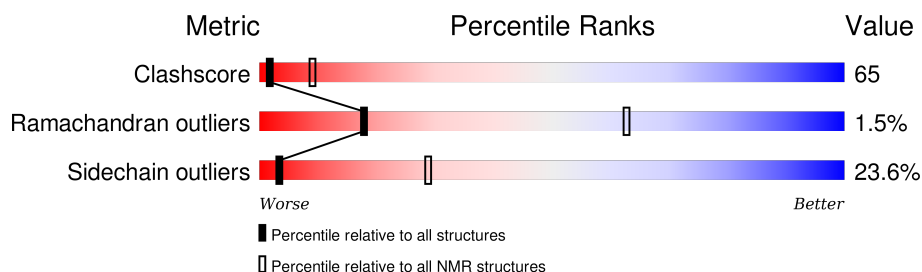
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 83%.

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)	NMR archive (#Entries)
Clashscore	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	109	

## 2 Ensemble composition and analysis ⓘ

This entry contains 23 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:122-A:222 (101)	0.27	7

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 4 clusters and 3 single-model clusters were found.

Cluster number	Models
1	1, 3, 4, 5, 6, 7, 11, 12, 14, 18, 21, 22
2	2, 16, 23
3	8, 15, 19
4	10, 20
Single-model clusters	9; 13; 17

### 3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1731 atoms, of which 877 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called HYPOTHETICAL 62.8 KDA PROTEIN C215.07C.

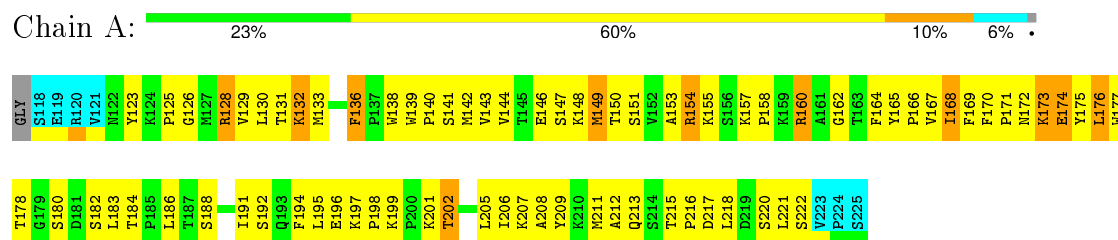
Mol	Chain	Residues	Atoms						Trace
1	A	108	Total	C	H	N	O	S	0
			1731	553	877	139	157	5	

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

### 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

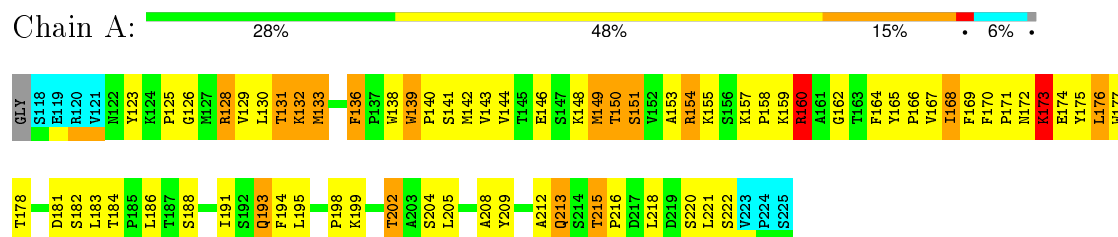
- Molecule 1: HYPOTHETICAL 62.8 KDA PROTEIN C215.07C



### 4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 7. Colouring as in section 4.1 above.

- Molecule 1: HYPOTHETICAL 62.8 KDA PROTEIN C215.07C



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *DISTANCE GEOMETRY SIMULATED ANNEALING*.

Of the 30 calculated structures, 23 were deposited, based on the following criterion: *NO VIOLATION GREATER THAN 0.25Å AND NO VIOLATIONS GREATER THAN 5 DEGREES*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
X-PLOR 3.8	refinement	
ANSIG 3.3	structure solution	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	BMRB entry 5538, BMRB entry 5902
Number of chemical shift lists	2
Total number of shifts	2495
Number of shifts mapped to atoms	1255
Number of unparsed shifts	0
Number of shifts with mapping errors	1240
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	83%

No validations of the models with respect to experimental NMR restraints is performed at this time.

## 6 Model quality [i](#)

### 6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	2.8±0.4
All	All	0	65

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	154	ARG	Sidechain	23
1	A	128	ARG	Sidechain	22
1	A	160	ARG	Sidechain	20

### 6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	800	821	821	105±9
All	All	18400	18883	18883	2426

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 65.

5 of 466 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:149:MET:HE3	1:A:153:ALA:HB1	1.02	1.27	4	10
1:A:215:THR:HG21	1:A:221:LEU:HD13	1.00	1.32	12	23
1:A:131:THR:CG2	1:A:178:THR:HG21	0.96	1.89	12	13
1:A:123:TYR:HB3	1:A:143:VAL:HG21	0.89	1.44	17	21
1:A:215:THR:CG2	1:A:221:LEU:HD13	0.88	1.97	9	19

## 6.3 Torsion angles [i](#)

### 6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	101/109 (93%)	93±2 (92±2%)	7±2 (7±2%)	2±1 (2±1%)	18	63
All	All	2323/2507 (93%)	2129 (92%)	158 (7%)	36 (2%)	18	63

All 4 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	150	THR	13
1	A	216	PRO	11
1	A	173	LYS	7
1	A	174	GLU	5

### 6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	91/98 (93%)	70±4 (76±4%)	21±4 (24±4%)	3	29
All	All	2093/2254 (93%)	1599 (76%)	494 (24%)	3	29

5 of 54 unique residues with a non-rotameric sidechain are listed below. They are sorted by the



frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	168	ILE	23
1	A	202	THR	23
1	A	173	LYS	22
1	A	136	PHE	21
1	A	132	LYS	18

### 6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 6.7 Other polymers [i](#)

There are no such molecules in this entry.

## 6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 83% for the well-defined parts and 83% for the entire structure.

### 7.1 Chemical shift list 1

File name: BMRB entry 5538

Chemical shift list name: *assigned\_chem\_shift\_list\_1*

#### 7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	1255
Number of shifts mapped to atoms	1255
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	7

#### 7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	107	$0.68 \pm 0.13$	Should be applied
$^{13}\text{C}_\beta$	99	$0.39 \pm 0.10$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}'$	107	$0.33 \pm 0.09$	None needed ( $< 0.5$ ppm)
$^{15}\text{N}$	97	$0.07 \pm 0.28$	None needed ( $< 0.5$ ppm)

#### 7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 83%, i.e. 1048 atoms were assigned a chemical shift out of a possible 1264. 13 out of 13 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	481/485 (99%)	192/192 (100%)	198/202 (98%)	91/91 (100%)
Sidechain	508/666 (76%)	374/400 (94%)	130/241 (54%)	4/25 (16%)

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	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Aromatic	59/113 (52%)	56/59 (95%)	0/51 (0%)	3/3 (100%)
Overall	1048/1264 (83%)	622/651 (96%)	328/494 (66%)	98/119 (82%)

#### 7.1.4 Statistically unusual chemical shifts [i](#)

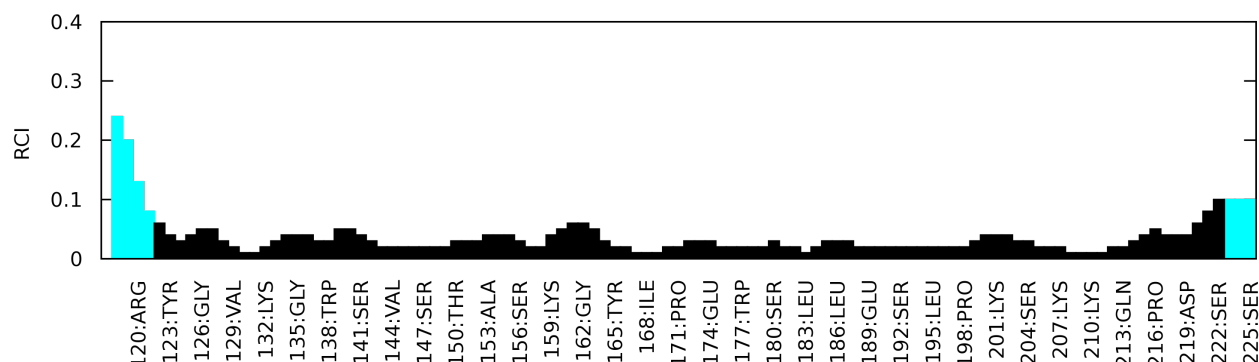
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	132	LYS	HG2	-1.16	2.67 – 0.07	-9.7
1	A	157	LYS	HG3	-0.80	2.76 – -0.04	-7.7
1	A	166	PRO	HB3	-0.39	3.81 – 0.21	-6.7
1	A	198	PRO	HA	2.40	6.05 – 2.75	-6.1
1	A	132	LYS	HG3	-0.23	2.76 – -0.04	-5.7
1	A	140	PRO	HG3	0.04	3.56 – 0.26	-5.7
1	A	157	LYS	HA	2.01	6.46 – 2.06	-5.1

#### 7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition.

Random coil index (RCI) for chain A:



## 7.2 Chemical shift list 2

File name: BMRB entry 5902

Chemical shift list name: *assigned\_chem\_shift\_list\_1*

### 7.2.1 Bookkeeping [i](#)

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	1240
Number of shifts mapped to atoms	0
Number of unparsed shifts	0
Number of shifts with mapping errors	1240
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	9

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- Chain not found in structure. First 5 (of 1240) occurrences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
UNMAPPED	106	LYS	HB3	1.56	0.02	2
UNMAPPED	65	THR	H	8.222	0.02	1
UNMAPPED	18	LYS	CE	42.155	0.2	1
UNMAPPED	73	LEU	CB	45.737	0.2	1
UNMAPPED	20	TYR	H	7.652	0.02	1

### 7.2.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	100	0.00 $\pm$ 0.00	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	92	0.00 $\pm$ 0.00	None needed (< 0.5 ppm)
$^{13}\text{C}'$	100	0.00 $\pm$ 0.00	None needed (< 0.5 ppm)
$^{15}\text{N}$	93	0.00 $\pm$ 0.00	None needed (< 0.5 ppm)

### 7.2.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 0%, i.e. 0 atoms were assigned a chemical shift out of a possible 1264. 0 out of 13 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	0/485 (0%)	0/192 (0%)	0/202 (0%)	0/91 (0%)
Sidechain	0/666 (0%)	0/400 (0%)	0/241 (0%)	0/25 (0%)
Aromatic	0/113 (0%)	0/59 (0%)	0/51 (0%)	0/3 (0%)
Overall	0/1264 (0%)	0/651 (0%)	0/494 (0%)	0/119 (0%)

### 7.2.4 Statistically unusual chemical shifts [i](#)

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
???	UNMAPPED	37	PRO	CD	42.04	55.31 – 45.41	-8.4
???	UNMAPPED	93	SER	CB	52.61	71.24 – 56.34	-7.5
???	UNMAPPED	103	PRO	CD	42.98	55.31 – 45.41	-7.5
???	UNMAPPED	101	ASN	HB3	0.42	4.41 – 1.11	-7.1
???	UNMAPPED	29	LYS	HD3	-0.01	2.75 – 0.45	-7.0
???	UNMAPPED	40	ILE	CG2	26.34	24.63 – 10.43	6.2
???	UNMAPPED	83	PHE	HB3	0.66	4.85 – 1.05	-6.0
???	UNMAPPED	87	ASN	HD21	4.90	9.74 – 4.94	-5.1
???	UNMAPPED	71	LYS	CG	19.16	30.67 – 19.17	-5.0

### 7.2.5 Random Coil Index (RCI) plots [i](#)

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned\_chem\_shift\_list\_1). RCI is only applicable to proteins.