

# Full wwPDB NMR Structure Validation Report (i)

## Oct 31, 2021 – 11:04 AM EDT

PDB ID	:	1P5L
Title	:	HP $(2-20)$ Substitution PHE5 to SER modification in sds-d25 micelles
Authors	:	Lee, K.H.; Lee, D.G.; Park, Y.K.; Harm, K.S.; Kim, Y.M.
Deposited on	:	2003-04-27

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org* A user guide is available at https://www.wwpdb.org/validation/2017/NMRValidationReportHelp with specific help available everywhere you see the (i) symbol.

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

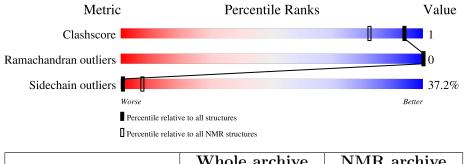
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. $(2010)$
ShiftChecker	:	2.23.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $SOLUTION\ NMR$ 

The overall completeness of chemical shifts assignment was not calculated.

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)	${f NMR}  ext{ archive} \ (\#  ext{Entries})$		
Clashscore	158937	12864		
Ramachandran outliers	154571	11451		
Sidechain outliers	154315	11428		

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 >=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 <=5%

Mol	Chain	Length	Quality of chain		
1	А	20	30%	15%	55%



# 2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 16 is the overall representative, medoid model (most similar to other models).

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:3-A:11 (9)	0.06	16		

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 1 single-model cluster was found.

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



# 3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 344 atoms, of which 185 are hydrogens and 0 are deuteriums.

• Molecule 1 is a protein called 19-mer peptide from 50S ribosomal protein L1.

Mol	Chain	Residues		At	$\mathbf{oms}$			Trace
1	А	20	Total 344	C 101	Н 185	N 31	0 27	1

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	5	SER	PHE	engineered mutation	UNP Q9ZK21



# 4 Residue-property plots (i)

## 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide 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: 19-mer peptide from 50S ribosomal protein L1

Chain	A:	30%	15%	55%	
A1 K2 K3	ЕЭ К10 L11	F12 S13 K14 115 Q16 N17 D18 K19 NH220			

## 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

## 4.2.1 Score per residue for model 1

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1



## 4.2.2 Score per residue for model 2





#### 4.2.3 Score per residue for model 3

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A:	25%	20%	55%	-
A1 K2 K3 R7 L8 E9 E9	L11 F12 F12 F12 I15 I15 Q16 Q16 D18 K19 M17 N17 D18			

#### 4.2.4 Score per residue for model 4

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain	A:	25%	20%	55%
A1 K2 K3 F9	K10 L11	F12 813 115 115 016 016 018 018 018 018		

#### 4.2.5 Score per residue for model 5

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

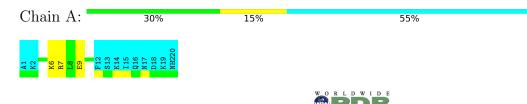
Chain A:	20%	25%	55%
A1 K2 K3 K6 R7	K10 L11 F12 S13 S13 115 115 115 N14 Y19 K19 K19	NH220	

#### 4.2.6 Score per residue for model 6

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A:	20%	20%	5%	55%
A1 K3 K7 K7	K10 E11 S13 S13 K14 T15 Q16 M16 M17 M17 X19	NH220		

#### 4.2.7 Score per residue for model 7



#### 4.2.8 Score per residue for model 8

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A:	30%	15%	55%	
A1 K3 K7 K10 K10	F11 F13 813 813 813 016 016 810 810 810 8120			

#### 4.2.9 Score per residue for model 9

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A	.:	20%		25%	55%	
A1 K2 K3 K6 K6 R7	L8 E9 K10 L11 F12	S13 K14 Q15 N17 D18	K19 NH220			

#### 4.2.10 Score per residue for model 10

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

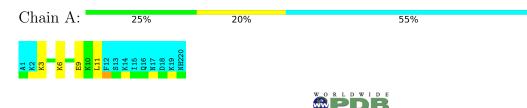
Chain A	.:	35%	10%	55%
A1 KG E9 E9	F12 513 513 115 115 016 013			

#### 4.2.11 Score per residue for model 11

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A:	30%	10% 5%	55%	-
A1 K3 S5 S5 K10	L11 F12 S13 F12 T15 Q16 Q16 D18 K19 NH220			

#### 4.2.12 Score per residue for model 12



#### 4.2.13 Score per residue for model 13

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A: 35% 10% 55%

### 4.2.14 Score per residue for model 14

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A: 35% 10% 55%

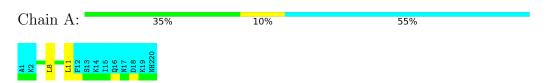
## 4.2.15 Score per residue for model 15

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

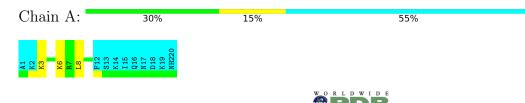
Chain A: 35% 10% 55%

## 4.2.16 Score per residue for model 16 (medoid)

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1



### 4.2.17 Score per residue for model 17



## 4.2.18 Score per residue for model 18

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A: 30% 15% 55%

#### 4.2.19 Score per residue for model 19

• Molecule 1: 19-mer peptide from 50S ribosomal protein L1

Chain A:	25%	20%	55%	
A1 K2 K3 K7 E9 K10	L11 F12 K14 K14 R15 N17 N17 K19 K19 N1220			

### 4.2.20 Score per residue for model 20

Chain A:	15%	25%	5%	55%
A1 K2 V4 K6 K6 K6 K6 K6 K6 K6 K6 K6 K6 K6 K6 K6	60 111 112 112 115 115 016	N17 D18 K19 NH220		



# 5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: HYBRID DISTANCE GEOMETRY - DYNAMICAL SIMULATED ANNEALING.

Of the 50 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
CNS	refinement	1.1

No chemical shift data was provided.



# 6 Model quality (i)

## 6.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: NH2

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

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

## 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	А	76	94	94	$0\pm 0$
All	All	1520	1880	1880	4

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 1.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
Atom-1	Atom-2	Clash(A)	Distance(A)	Worst	Total
1:A:11:LEU:HD12	1:A:11:LEU:C	0.45	2.31	12	1
1:A:8:LEU:HA	1:A:11:LEU:HD12	0.41	1.92	2	1
1:A:3:LYS:C	1:A:5:SER:N	0.41	2.74	20	2

All unique clashes are listed below, sorted by their clash magnitude.

## 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



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	А	9/20~(45%)	8±0 (86±5%)	$1\pm0~(14\pm5\%)$	0±0 (0±0%)	100 100
All	All	180/400~(45%)	155 (86%)	25 (14%)	0 (0%)	100 100

was analysed and the total number of residues.

There are no Ramachandran outliers.

### 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	А	9/18~(50%)	$6\pm1~(63\pm12\%)$	$3\pm1$ (37 $\pm12\%$ )	1 7
All	All	180/360~(50%)	113 (63%)	67~(37%)	1 7

All 7 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	А	3	LYS	14
1	А	10	LYS	14
1	А	9	GLU	11
1	А	7	ARG	10
1	А	6	LYS	9
1	А	11	LEU	6
1	А	8	LEU	3

## 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 monosaccharides 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 (i)

No chemical shift data were provided

