



# wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 2L22  
Title : Mupirocin didomain ACP  
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Deposited on : 2010-08-10

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  
<https://www.wwpdb.org/validation/2017/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  
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 : 2.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

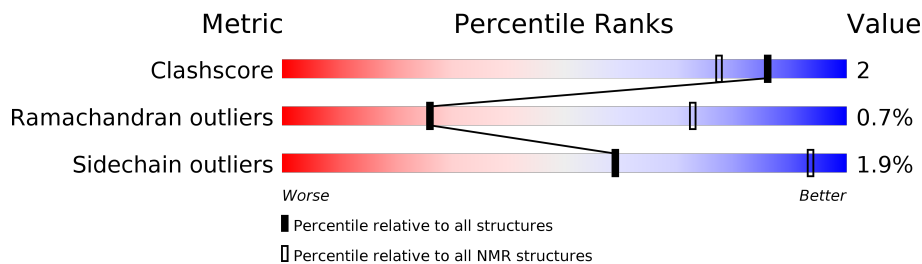
# 1 Overall quality at a glance i

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)	NMR archive (#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  $\geq 3$ , 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	212	

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 5 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

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:4-A:78 (75)	0.65	9
2	A:104-A:183 (80)	0.99	5

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 2 clusters. No single-model clusters were found.

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

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

There is only 1 type of molecule in this entry. The entry contains 2850 atoms, of which 1414 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Mupirocin didomain Acyl Carrier Protein.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	183	2850	910	1414	250	271	5	0

There are 29 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	EXPRESSION TAG	UNP Q8RL76
A	-19	GLY	-	EXPRESSION TAG	UNP Q8RL76
A	-18	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-17	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-16	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-15	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-14	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-13	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-12	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-11	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-10	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-9	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-8	GLY	-	EXPRESSION TAG	UNP Q8RL76
A	-7	LEU	-	EXPRESSION TAG	UNP Q8RL76
A	-6	VAL	-	EXPRESSION TAG	UNP Q8RL76
A	-5	PRO	-	EXPRESSION TAG	UNP Q8RL76
A	-4	ARG	-	EXPRESSION TAG	UNP Q8RL76
A	-3	GLY	-	EXPRESSION TAG	UNP Q8RL76
A	-2	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-1	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	0	MET	-	EXPRESSION TAG	UNP Q8RL76
A	184	LEU	-	EXPRESSION TAG	UNP Q8RL76
A	185	GLU	-	EXPRESSION TAG	UNP Q8RL76
A	186	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	187	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	188	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	189	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	190	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	191	HIS	-	EXPRESSION TAG	UNP Q8RL76



## 5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *simulated annealing*.

Of the 100 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
ARIA	refinement	2.2

No chemical shift data was provided. No validations of the models with respect to experimental NMR restraints is performed at this time.

COVALENT-GEOMETRY INFOmissingINFO

### 5.1 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	1242	1225	1223	5±2
All	All	24840	24500	24460	103

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.

5 of 70 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:164:ALA:HA	1:A:168:TRP:CZ3	0.62	2.29	16	1
1:A:110:ILE:HG21	1:A:178:VAL:HG11	0.60	1.73	3	6
1:A:60:GLY:HA2	1:A:73:TRP:CZ3	0.60	2.32	7	1
1:A:30:ARG:HA	1:A:65:PRO:O	0.59	1.97	8	12
1:A:27:ASP:HA	1:A:67:PHE:CZ	0.56	2.36	13	2

## 5.2 Torsion angles [i](#)

### 5.2.1 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	154/212 (73%)	146±2 (95±1%)	7±2 (5±1%)	1±1 (1±1%)	26	73
All	All	3080/4240 (73%)	2914 (95%)	144 (5%)	22 (1%)	26	73

5 of 14 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	123	PHE	4
1	A	129	LEU	3
1	A	130	ARG	2
1	A	157	SER	2
1	A	132	GLY	2

### 5.2.2 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	131/176 (74%)	129±1 (98±1%)	2±1 (2±1%)	59	93
All	All	2620/3520 (74%)	2571 (98%)	49 (2%)	59	93

5 of 34 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	149	MET	4
1	A	5	GLU	3
1	A	134	GLN	3
1	A	177	GLU	3
1	A	47	GLU	2

### 5.2.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.4 Carbohydrates [i](#)

There are no carbohydrates in this entry.

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

There are no ligands in this entry.

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

There are no such molecules in this entry.

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

There are no chain breaks in this entry.



## 6 Chemical shift validation

No chemical shift data were provided