

Full wwPDB NMR Structure Validation Report (i)

Mar 1, 2022 – 10:52 AM EST

PDB ID	:	2G9L
Title	:	The High-resolution Solution Conformation of an Antimicrobial Peptide Gae-
		gurin 4 and Its Mode of Membrane Interaction
Authors	:	Chi, SW.; Han, KH.
Deposited on	:	2006-03-06

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 (1)) were used in the production of this report:

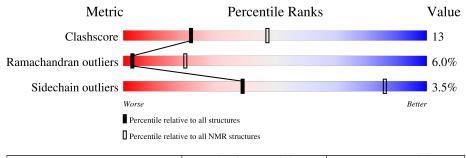
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.27
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.27

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 $(\#Fntricg)$	${f NMR} \; { m archive} \ (\#{ m Entries})$			
	$(\# { m Entries})$	(#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 c	hain	
1	А	37	57%	30%	14%



2 Ensemble composition and analysis (i)

This entry contains 15 models. Model 10 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:3-A:22 (20)	0.30	10				
2	A:25-A:36 (12)	0.49	9				

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. No single-model clusters were found.

Cluster number	Models
1	2, 4, 6, 8, 10, 11
2	1, 5, 7, 9, 14
3	3, 13
4	12, 15



3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 553 atoms, of which 292 are hydrogens and 0 are deuteriums.

• Molecule 1 is a protein called Gaegurin-4.

Mol	Chain	Residues	Atoms					Trace	
1	٨	27	Total	С	Η	Ν	Ο	S	0
	A	51	553	165	292	45	49	2	0

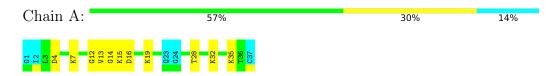


4 Residue-property plots (i)

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.

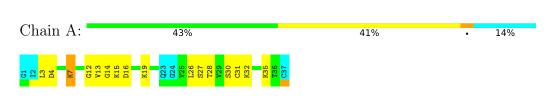
• Molecule 1: Gaegurin-4



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



4.2.2 Score per residue for model 2

 \bullet Molecule 1: Gaegurin-4





4.2.3 Score per residue for model 3

• Molecule 1: Gaegurin-4



4.2.4 Score per residue for model 4

• Molecule 1: Gaegurin-4

Chain A:				709	%			16%	14%
G1 12 V13 K15 D16	K1 <mark>9</mark> Q23 G24	T28 K32	<mark>C37</mark>						

4.2.5 Score per residue for model 5

• Molecule 1: Gaegurin-4

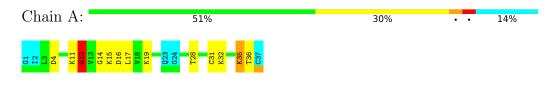
Chain A:	62%	22%	• 14%
G1 12 12 12 12 12 12 12 12 12 14 13 16 116 116 116	023 024 024 037 037 037		

4.2.6 Score per residue for model 6

• Molecule 1: Gaegurin-4



4.2.7 Score per residue for model 7





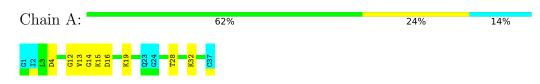
4.2.8 Score per residue for model 8

• Molecule 1: Gaegurin-4



4.2.9 Score per residue for model 9

• Molecule 1: Gaegurin-4



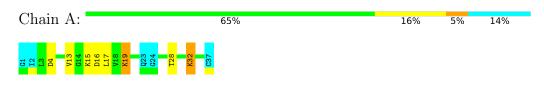
4.2.10 Score per residue for model 10 (medoid)

• Molecule 1: Gaegurin-4

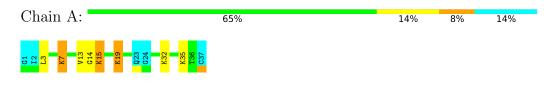
Chain A:	62%	24%	14%
61 12 13 14 14 14 14 15 14 116 116 116 116 116 116 116 116 116	128 K32 K35 C37 C37		

4.2.11 Score per residue for model 11

• Molecule 1: Gaegurin-4



4.2.12 Score per residue for model 12





4.2.13 Score per residue for model 13

• Molecule 1: Gaegurin-4



4.2.14 Score per residue for model 14

• Molecule 1: Gaegurin-4



4.2.15 Score per residue for model 15

Chain A:	65%	19%	·	14%
01 12 13 14 13 14 13 14 13 14 13 12 13 13 13 13 13 13 13 13 13 13 13 13 13				



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: distance geometry, molecular dynamics.

Of the 50 calculated structures, 15 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
Discover	refinement	2.98

No chemical shift data was provided.



6 Model quality (i)

6.1 Standard geometry (i)

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 (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mal	Mol Chain		ond lengths	I	Bond angles
	Unam	RMSZ	#Z > 5	RMSZ	$\#Z{>}5$
1	А	$0.63 {\pm} 0.00$	$2{\pm}1/2{30}~(~0.8{\pm}~0.2\%)$	$0.68 {\pm} 0.02$	$0{\pm}0/308~(~0.0{\pm}~0.0\%)$
All	All	0.63	26/3450 ($0.8%$)	0.68	0/4620 ($0.0%$)

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	А	$0.0{\pm}0.0$	0.3 ± 0.4
All	All	0	4

All unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Dec	Turne	Atoms	7	Observed(Å)	$I_{doal}(\lambda)$	Moo	dels
IVIOI	Unam	nes	Type	Atoms		Observed(A)	Ideal(A)	Worst	Total
1	А	4	ASP	CG-OD2	5.05	1.36	1.25	15	12
1	А	16	ASP	CG-OD2	5.05	1.36	1.25	7	14

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	А	12	GLY	Peptide	2
1	А	28	THR	Peptide	1
1	А	32	LYS	Peptide	1



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	229	260	258	6 ± 3
All	All	3435	3900	3870	93

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 13.

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

Atom 1	Atom 2	$Clash(\lambda)$	Distance (Å)	Moo	dels
Atom-1	Atom-2	Clash(Å)	$\operatorname{Distance}(\operatorname{\AA})$	Worst	Total
1:A:32:LYS:HD3	1:A:35:LYS:CD	1.33	1.53	5	2
1:A:32:LYS:CD	1:A:35:LYS:HD3	1.30	1.57	14	2
1:A:32:LYS:CD	1:A:35:LYS:CD	1.24	2.15	5	2
1:A:31:CYS:HB2	1:A:35:LYS:CD	1.23	1.63	7	1
1:A:31:CYS:CB	1:A:35:LYS:HD2	1.17	1.68	7	1
1:A:32:LYS:CD	1:A:35:LYS:HD2	1.13	1.71	5	1
1:A:3:LEU:O	1:A:7:LYS:HD3	1.07	1.50	8	2
1:A:32:LYS:HD2	1:A:35:LYS:HD3	1.06	1.11	14	1
1:A:3:LEU:O	1:A:7:LYS:HD2	1.02	1.52	12	2
1:A:32:LYS:HD3	1:A:35:LYS:HD3	1.01	1.31	5	1
1:A:32:LYS:HD2	1:A:35:LYS:HE3	1.00	1.28	5	2
1:A:28:THR:O	1:A:32:LYS:HB3	0.96	1.59	2	2
1:A:32:LYS:HD3	1:A:35:LYS:HD2	0.96	1.20	5	2
1:A:32:LYS:HD2	1:A:35:LYS:CE	0.95	1.91	5	2
1:A:15:LYS:O	1:A:19:LYS:HB2	0.92	1.65	3	8
1:A:3:LEU:O	1:A:7:LYS:CD	0.91	2.18	12	3
1:A:28:THR:O	1:A:32:LYS:HB2	0.90	1.67	15	9
1:A:27:SER:O	1:A:31:CYS:CB	0.87	2.22	1	1
1:A:15:LYS:O	1:A:19:LYS:CB	0.83	2.27	6	7
1:A:28:THR:O	1:A:32:LYS:HD3	0.78	1.77	6	1
1:A:27:SER:O	1:A:31:CYS:HB2	0.76	1.78	1	1
1:A:32:LYS:CD	1:A:35:LYS:CE	0.74	2.61	5	2
1:A:32:LYS:CE	1:A:35:LYS:HD2	0.73	2.14	5	1
1:A:28:THR:O	1:A:32:LYS:CD	0.67	2.42	6	1
1:A:27:SER:O	1:A:31:CYS:HB3	0.66	1.88	1	1
1:A:31:CYS:HB3	1:A:36:THR:HB	0.61	1.72	8	1
1:A:28:THR:HG22	1:A:32:LYS:HD2	0.58	1.75	6	1

Continued on next page...



Atom-1	Atom-2	Clash(Å)	Distance(Å)	Moo	dels
Atom-1	Atom-2	Clash(A)	Distance(A)	Worst	Total
1:A:31:CYS:HB2	1:A:35:LYS:HD2	0.58	0.74	7	1
1:A:12:GLY:O	1:A:16:ASP:OD2	0.57	2.21	14	1
1:A:28:THR:O	1:A:32:LYS:CB	0.55	2.46	2	5
1:A:3:LEU:O	1:A:7:LYS:HB2	0.54	2.02	2	1
1:A:26:LEU:O	1:A:30:SER:HB3	0.54	2.03	1	1
1:A:32:LYS:O	1:A:34:ALA:N	0.53	2.42	6	1
1:A:32:LYS:O	1:A:32:LYS:HD3	0.52	2.05	11	1
1:A:26:LEU:O	1:A:30:SER:CB	0.50	2.60	1	1
1:A:19:LYS:HD3	1:A:19:LYS:C	0.50	2.27	6	3
1:A:15:LYS:O	1:A:19:LYS:N	0.49	2.44	12	7
1:A:7:LYS:N	1:A:7:LYS:CD	0.48	2.77	10	1
1:A:12:GLY:H	1:A:14:GLY:H	0.48	1.52	3	1
1:A:11:LYS:O	1:A:15:LYS:HE3	0.47	2.09	7	1
1:A:12:GLY:C	1:A:14:GLY:H	0.46	2.13	5	2
1:A:7:LYS:HE3	1:A:11:LYS:HD2	0.45	1.89	6	1
1:A:12:GLY:N	1:A:14:GLY:H	0.43	2.12	7	2
1:A:32:LYS:C	1:A:32:LYS:HD3	0.41	2.37	9	1
1:A:32:LYS:CD	1:A:35:LYS:HE2	0.40	2.45	12	1
1:A:19:LYS:HD3	1:A:19:LYS:O	0.40	2.15	6	1

Continued from previous page...

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	Pe	erce	entiles
1	А	32/37~(86%)	26 ± 2 (81 $\pm5\%$)	$4\pm1~(13\pm4\%)$	$2\pm1~(6\pm3\%)$		3	20
All	All	480/555~(86%)	389 (81%)	62 (13%)	29~(6%)		3	20

All 5 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	А	13	VAL	13
1	А	12	GLY	8
1	А	14	GLY	5

Continued on next page...



Continued from previous page...

Mol	Chain	Res	Type	Models (Total)
1	А	33	LEU	2
1	А	35	LYS	1

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	25/28~(89%)	$24 \pm 1 (97 \pm 4\%)$	$1\pm1 (3\pm4\%)$	39 86
All	All	375/420 (89%)	362 (97%)	13 (3%)	39 86

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	32	LYS	4
1	А	7	LYS	2
1	А	17	LEU	2
1	А	19	LYS	2
1	А	13	VAL	1
1	А	35	LYS	1
1	А	15	LYS	1

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

