

# Full wwPDB NMR Structure Validation Report (i)

### Sep 7, 2022 – 07:55 PM EDT

PDB ID	:	7UZL
Title	:	Solution NMR structure of 9-residue Rosetta-designed cyclic peptide D9.16 in
		CDCl3 with cis/trans switching (B-TC conformation)
Authors	:	Ramelot, T.A.; Tejero, R.; Montelione, G.T.
Deposited on	:	2022-05-09

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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

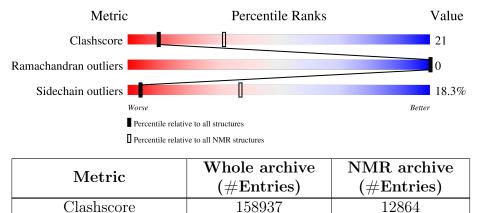
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	$v_1n_11_5_13_A$ (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.29
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.29

#### Overall quality at a glance (i) 1

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

The overall completeness of chemical shifts assignment is 94%.

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.



Ramachandran outliers	154571	11451			
Sidechain outliers	154315	11428			
			-		
			1		
	0		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 fract	tion of residues that a	re not part of the we	ll-defined cores, and a grey seg-		

ment 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	Qua	lity of chain	
1	А	9	44%	44%	11%



# 2 Ensemble composition and analysis (i)

This entry contains 20 models.

Cyrange was unable to find well-defined residues.

Error message: The number of core atoms (4) was below the domain threshold value (8).

NmrClust was unable to cluster the ensemble.

Error message: Wrapper check: not enough residues in core to run NmrClust



# 3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 146 atoms, of which 81 are hydrogens and 0 are deuteriums.

• Molecule 1 is a protein called Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU -PRO-DLE.

[	Mol	Chain	Residues		Ate	$\mathbf{oms}$			Trace
	1	۸	0	Total	С	Η	Ν	0	0
	1	A	9	146	47	81	9	9	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: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	44%	44%	11%
P1 L5 L6 L7 L7 L7 L9 L9 L9			

## 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: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	56%	44%	•
P1 L5 L7 L6 L6 L9			

## 4.2.2 Score per residue for model 2

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A: 67% 22% 11%



#### 4.2.3 Score per residue for model 3

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	44%	44%	11%
<mark>г5</mark> г5 г3 г3 г3			

#### 4.2.4 Score per residue for model 4

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	44%	44%	11%
12 12 12 12 12 12 12 12 12 12 12 12 12 1			

#### 4.2.5 Score per residue for model 5

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	44%	56%	
P1 L5 L6 P8 L9 L9			

#### 4.2.6 Score per residue for model 6

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	56%	33%	11%
2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2			

#### 4.2.7 Score per residue for model 7

Chain A:	44%	56%
1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1		



#### 4.2.8 Score per residue for model 8

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	44%	56%	-
L L P B B B B B B B C C C C C C C C C C C C			

#### 4.2.9 Score per residue for model 9

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	44%	33%	22%
<mark>1</mark> 15 15 15 15 15 15 15 15 15 15 15 15 15			

#### 4.2.10 Score per residue for model 10

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	44%	33%	22%
10 10 10 10 10 10 10 10 10 10 10 10 10 1			

#### 4.2.11 Score per residue for model 11

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	33%	56%	11%	
P1 L5 L5 P8 L9 L9				

#### 4.2.12 Score per residue for model 12

Chain A:	44%	56%
<b>1</b> 12 12 13 13 14 14 14 14 14 14 14 14 14 14 14 14 14		



#### 4.2.13 Score per residue for model 13

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	44%	44%	11%
L1 L1 L2 L3 L3 L3 L3 L3 L3 L3 L3 L3 L3 L3 L3 L3			

4.2.14 Score per residue for model 14

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	67%	33%	•
14 91 17 17 17 17 17 17 17 17 17 17 17 17 17			

#### 4.2.15 Score per residue for model 15

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

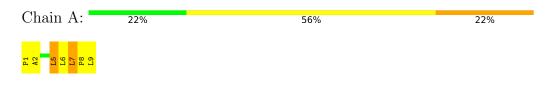
Chain A:	44%	44%	11%
P1 L5 L7 P8 L9			

#### 4.2.16 Score per residue for model 16

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A: 44% 11% 로 입입되었었

#### 4.2.17 Score per residue for model 17



#### 4.2.18 Score per residue for model 18

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A: 11% 78% 11%

P1 A2 A3 L5 L5 L6 L6 L6 L7 L7 L7 L9

#### 4.2.19 Score per residue for model 19

• Molecule 1: Cyclic peptide D9.16 DPR-MAA-ALA-DVA-MLE-LEU-LEU-PRO-DLE

Chain A:	22%	56%	22%
P1 L5 L5 L6 L7 P8 P8			

#### 4.2.20 Score per residue for model 20

Chain A:	56%	33%	11%
2000 100 100 100 100 100 100 100 100 100			



# 5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: na.

Of the 100 calculated structures, 20 were deposited, based on the following criterion: target function.

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

Software name	Classification	Version
CYANA	refinement	3.98.13
CYANA	structure calculation	3.98.13

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)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	112
Number of shifts mapped to atoms	112
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	94%



# 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: MLE, DPR, MAA, DLE, DVA

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	65	81	80	$3\pm1$
All	All	1300	1620	1598	62

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 21.

Models Atom-1 Atom-2 Clash(Å) Distance(Å)Worst Total 1:A:4:DVA:HG22 1:A:5:MLE:HD13 0.811.52191 1:A:6:LEU:C 1:A:9:DLE:HD22 0.642.13127 1:A:4:DVA:HG22 1:A:5:MLE:HG 0.551.7911 1 1 1:A:4:DVA:HG22 1:A:5:MLE:CD1 0.532.29191:A:5:MLE:HN3 1:A:6:LEU:HD23 203 0.531.81 1:A:5:MLE:HN3 1:A:6:LEU:HD12 0.529 3 1.81 1:A:4:DVA:HG12 1:A:5:MLE:HG 1 0.521.81 18 1:A:6:LEU:N 1:A:9:DLE:HD22 0.512.201913 1:A:7:LEU:N 1:A:9:DLE:HD22 0.462.2612 1 2.43 1:A:1:DPR:CD 1:A:9:DLE:HD13 191 0.431:A:7:LEU:HD23 1:A:7:LEU:O 0.432.14 $\mathbf{2}$ 6 1:A:7:LEU:HA 1:A:8:PRO:C  $2.3\overline{4}$ 200.4316

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

Continued on next page...



Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
Atom-1	Atom-2	Clash(A)	Distance(A)	Worst	Total
1:A:6:LEU:HD12	1:A:7:LEU:O	0.43	2.13	13	1
1:A:1:DPR:HA	1:A:2:MAA:HM1	0.42	1.77	17	2
1:A:5:MLE:HN3	1:A:6:LEU:N	0.41	2.29	10	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	Percentile	es
1	А	4/9~(44%)	$2\pm0~(51\pm5\%)$	$2\pm0~(49\pm5\%)$	0±0 (0±0%)	100 100	)
All	All	80/180 (44%)	41 (51%)	39~(49%)	0  (0%)	100 100	)

There are no Ramachandran outliers.

#### 6.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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 side chain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
1	А	3/3~(100%)	$2\pm0$ (82 $\pm17\%$ )	$1\pm0 (18\pm17\%)$	4	37	
All	All	60/60~(100%)	49 (82%)	11 (18%)	4	37	

All 2 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	7	LEU	10
1	А	6	LEU	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)

5 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with |Z| > 2 is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Turne	Chain	Dec	Tink	Bond leng		gths
	туре	Unam	nes		Counts	RMSZ	#Z>2
1	MLE	А	5	1	$7,\!8,\!9$	$0.44{\pm}0.01$	0±0 (0±0%)
1	MAA	А	2	1	4,5,6	$0.55 {\pm} 0.01$	0±0 (0±0%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with |Z| > 2 is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mal	Trune	Chain	Dec	Timle	Bond angles Counts   RMSZ   #Z>2		
INIOI	туре	Chain	nes	LIIIK	Counts	RMSZ	$\#Z{>}2$
1	MLE	А	5	1	6,9,11	$1.48 {\pm} 0.00$	$2\pm0$ (33 $\pm0\%$ )
1	MAA	А	2	1	$1,\!5,\!7$	$0.30 {\pm} 0.01$	0±0 (0±0%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLE	А	5	1	-	$0\pm 0,5,8,10$	-
1	MAA	А	2	1	-	$0\pm 0,1,4,6$	-

There are no bond-length outliers.



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

Mol			$Ideal(^{o})$	Moo					
	Ullalli	Ites	Type	Atoms			Iueai()	Worst	Total
1	А	5	MLE	CG-CB-CA	2.37	109.43	115.34	14	20
1	А	5	MLE	CN-N-CA	2.24	106.66	113.64	18	20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

# 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)

The completeness of assignment taking into account all chemical shift lists is 94% for the well-defined parts and 94% for the entire structure.

# 7.1 Chemical shift list 1

File name: working\_cs.cif

Chemical shift list name: *chem\_shift\_list\_1* 

## 7.1.1 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	112
Number of shifts mapped to atoms	112
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

## 7.1.2 Chemical shift referencing (i)

No chemical shift referencing corrections were calculated (not enough data).

### 7.1.3 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 94%, i.e. 44 atoms were assigned a chemical shift out of a possible 47. 0 out of 2 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^{1}\mathbf{H}$	$^{13}\mathbf{C}$	$^{15}\mathbf{N}$
Backbone	15/18~(83%)	7/7~(100%)	6/8~(75%)	2/3~(67%)
Sidechain	29/29~(100%)	17/17~(100%)	12/12~(100%)	0/0 (%)
Aromatic	0/0~(-%)	0/0 (%)	0/0~(-%)	0/0 (%)
Overall	44/47~(94%)	24/24~(100%)	18/20~(90%)	2/3~(67%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 94%, i.e. 44 atoms were assigned a chemical shift out of a possible 47. 0 out of 2 assigned methyl groups (LEU and VAL) were assigned stereospecifically.



	Total	$^{1}\mathrm{H}$	$^{13}\mathrm{C}$	$^{15}\mathbf{N}$
Backbone	15/18~(83%)	7/7~(100%)	6/8~(75%)	2/3~(67%)
Sidechain	29/29~(100%)	17/17~(100%)	12/12~(100%)	$0/0 \ (\%)$
Aromatic	0/0~(-%)	0/0~(-%)	0/0~(-%)	0/0 (%)
Overall	44/47~(94%)	24/24~(100%)	18/20~(90%)	2/3~(67%)

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

There are no statistically unusual chemical shifts.

### 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:



