wwPDB NMR Structure Validation Summary Report (i)

Apr 26, 2016 – 02:28 PM BST

PDB ID	:	1DLZ
Title	:	SOLUTION STRUCTURE OF THE CHANNEL-FORMER ZERVAMICIN IIB (PEPTAIBOL ANTIBIOTIC)
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Deposited on	:	1999-12-13

This is a wwPDB NMR Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/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:

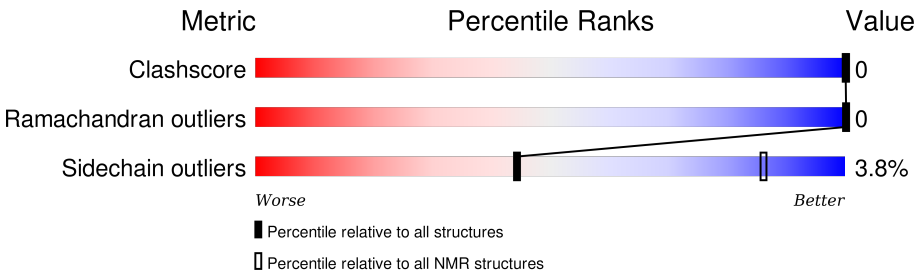
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	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 (i)

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

The overall completeness of chemical shifts assignment is 98%.

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 >=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	A	17	94% 6%

2 Ensemble composition and analysis (i)

This entry contains 20 models.

Cyrange was unable to find well-defined residues.

Error message: Only domains with < 8 residues could be identified.

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 270 atoms, of which 139 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called ZERVAMICIN IIB.

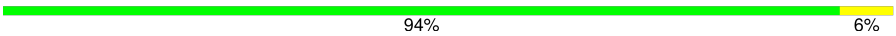
Mol	Chain	Residues	Atoms					Trace
			Total	C	H	N	O	
1	A	17	270	90	139	19	22	0

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: ZERVAMICIN IIB

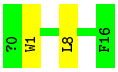
Chain A: 

4.2 Residue scores for the first model from the NMR ensemble

No representative models were identified. Colouring as in section 4.1 above.

- Molecule 1: ZERVAMICIN IIB

Chain A:



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: *BEST 20 OUT OF 100 WITH CA BACKBONE RMSD NO LESS THAN 0.05 ANGSTROMS*.

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

Software name	Classification	Version
DISCOVER 3.1	refinement	
DYANA 1.5	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 4601
Number of chemical shift lists	1
Total number of shifts	209
Number of shifts mapped to atoms	192
Number of unparsed shifts	0
Number of shifts with mapping errors	17
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	98%

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)

Bond lengths and bond angles in the following residue types are not validated in this section: HYP, DIV, PHL, ACE, AIB

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).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	1.21±0.01	0±0/71 (0.0±0.0%)	1.34±0.05	1±0/95 (1.1±0.0%)
All	All	1.21	0/1420 (0.0%)	1.34	20/1900 (1.1%)

There are no bond-length outliers.

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})	Models	
								Worst	Total
1	A	1	TRP	CD1-NE1-CE2	-5.71	103.86	109.00	17	20

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
All	All	2620	2780	2740	-

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 -.

There are no clashes.

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	8/17 (47%)	7±0 (88±0%)	1±0 (12±0%)	0±0 (0±0%)	100	100
All	All	160/340 (47%)	140 (88%)	20 (12%)	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 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	8/8 (100%)	8±0 (96±6%)	0±0 (4±6%)	44	86
All	All	160/160 (100%)	154 (96%)	6 (4%)	44	86

All 4 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	11	GLN	2
1	A	2	ILE	2
1	A	8	LEU	1
1	A	5	ILE	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)

8 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	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
1	HYP	A	10	1	6,8,9	0.43±0.01	0±0 (0±0%)
1	AIB	A	12	1	2,5,6	1.19±0.02	0±0 (0±0%)
1	HYP	A	13	1	6,8,9	0.44±0.01	0±0 (0±0%)
1	AIB	A	14	1	2,5,6	1.11±0.01	0±0 (0±0%)
1	PHL	A	16	1	11,11,11	0.97±0.01	0±0 (0±0%)
1	DIV	A	4	1	3,6,7	1.03±0.03	0±0 (0±0%)
1	AIB	A	7	1	2,5,6	1.15±0.02	0±0 (0±0%)
1	AIB	A	9	1	2,5,6	1.20±0.02	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.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
1	HYP	A	10	1	5,10,12	1.34±0.08	0±0 (0±0%)
1	AIB	A	12	1	4,7,9	3.37±0.06	1±1 (12±12%)
1	HYP	A	13	1	5,10,12	1.33±0.04	0±0 (0±0%)
1	AIB	A	14	1	4,7,9	2.93±0.06	0±0 (0±0%)
1	PHL	A	16	1	10,13,13	0.49±0.06	0±0 (0±0%)
1	DIV	A	4	1	3,8,10	1.70±0.07	0±0 (0±0%)
1	AIB	A	7	1	4,7,9	1.90±0.10	0±0 (0±0%)
1	AIB	A	9	1	4,7,9	2.97±0.05	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	HYP	A	10	1	-	0±0,0,11,13	0±0,1,1,1
1	AIB	A	12	1	-	0±0,2,3,6	0±0,0,0,0
1	HYP	A	13	1	-	0±0,0,11,13	0±0,1,1,1
1	AIB	A	14	1	-	0±0,2,3,6	0±0,0,0,0
1	PHL	A	16	1	-	0±0,6,6,6	0±0,1,1,1
1	DIV	A	4	1	-	0±0,3,6,9	0±0,0,0,0
1	AIB	A	7	1	-	0±0,2,3,6	0±0,0,0,0
1	AIB	A	9	1	-	0±0,2,3,6	0±0,0,0,0

There are no bond-length outliers.

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})	Models	
								Worst	Total
1	A	12	AIB	CB2-CA-CB1	5.29	102.31	110.42	7	10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 (i)

The completeness of assignment taking into account all chemical shift lists is 98% for the well-defined parts and 98% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 4601

Chemical shift list name: *assigned_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	209
Number of shifts mapped to atoms	192
Number of unparsed shifts	0
Number of shifts with mapping errors	17
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

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

- No matching atoms in chemical component dictionary. First 5 (of 17) occurences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
A	4	DIV	CG	8.91	0.1	1
A	16	PHL	CE	130.26	0.1	2
A	14	AIB	HB1	1.54	0.002	1
A	12	AIB	HB2	1.51	0.002	1
A	9	AIB	HB2	1.53	0.002	1

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	11	-0.48 ± 0.42	None needed (< 0.5 ppm)
$^{13}C_{\beta}$	11	0.00 ± 0.00	None needed (< 0.5 ppm)

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Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C'$	16	0.77 ± 0.24	Should be applied
^{15}N	13	0.45 ± 1.98	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments (i)

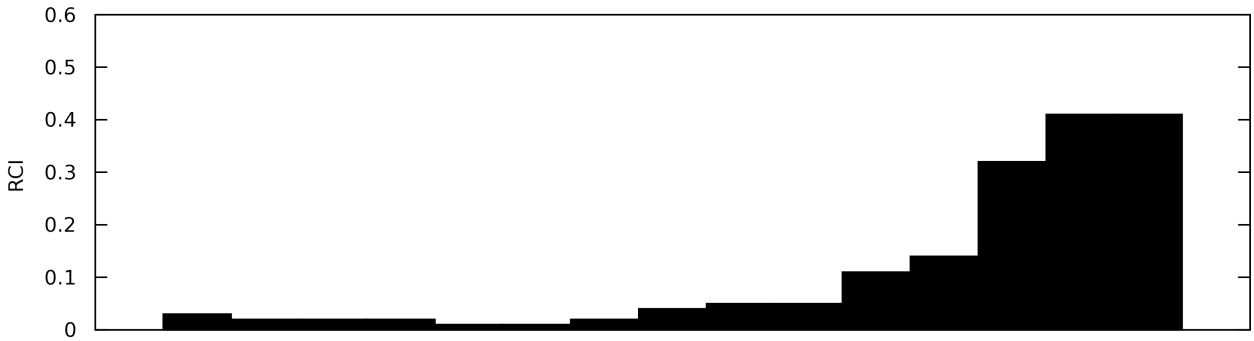
The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 98%, i.e. 111 atoms were assigned a chemical shift out of a possible 113. 0 out of 1 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	38/38 (100%)	15/15 (100%)	16/16 (100%)	7/7 (100%)
Sidechain	61/63 (97%)	36/37 (97%)	23/24 (96%)	2/2 (100%)
Aromatic	12/12 (100%)	6/6 (100%)	5/5 (100%)	1/1 (100%)
Overall	111/113 (98%)	57/58 (98%)	44/45 (98%)	10/10 (100%)

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: