wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1EWW
Title	:	SOLUTION STRUCTURE OF SPRUCE BUDWORM ANTIFREEZE PROTEIN AT 30 DEGREES CELSIUS
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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
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:

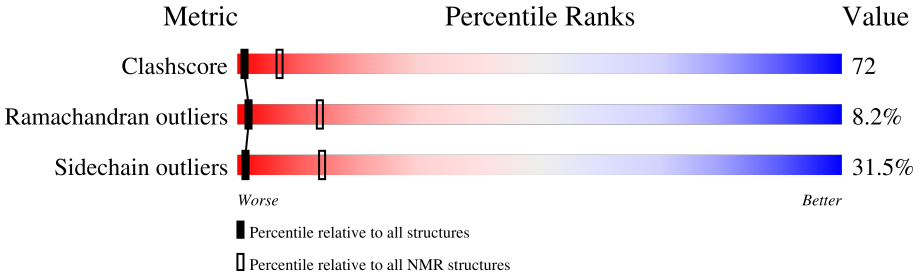
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.26
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.26

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 >=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	90	6% 61% 14% • 18%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 17 is the overall representative, medoid model (most similar to other models). The authors have identified model 10 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:18, A:24-A:72, A:81-A:90 (74)	0.38	17

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 and 3 single-model clusters were found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1229 atoms, of which 603 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called ANTIFREEZE PROTEIN.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	90	1229	372	603	107	139	8	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: ANTIFREEZE PROTEIN

Chain A: 6% 61% 14% • 18%

D1 G2 S3 C4 T5 N6 T7 N8 S9 Q10 L11 S12 A13 N14 S15 K16 C17 E18 K19 S20 T21 L22 T23 N24 C25 Y26 V27 D28 K29 S30 E31 V32 Y33 G34 T35 T36 C37 T38 G39 S40 R41 F42 D43 G44 V45 T46 I47 T48 T49 S50 T51 S52 T53 G54 S55 R56 I57 S58 G59

C62 K63 I64 S65 T66 C67 I68 I69 T70 G71 G72 V73 P74 A75 P76 S77 A78 A79 C80 K81 I82 S83 G84 C85 T86 F87 S88 A89 N90

4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 17. Colouring as in section 4.1 above.

• Molecule 1: ANTIFREEZE PROTEIN

Chain A: 16% 46% 20% • 18%

D1 G2 S3 C4 T5 N6 T7 N8 S9 Q10 L11 S12 A13 N14 S15 K16 C17 E18 K19 S20 T21 L22 T23 N24 C25 Y26 V27 D28 K29 S30 E31 V32 Y33 G34 T35 T36 C37 S40 R41 F42 D43 G44 V45 T46 I47 T48 T49 S50 T51 S52 T53 G54 S55 R56 I57 S58 G59 C62

K63 I64 S65 T66 C67 T70 G71 G72 V73 P74 A75 P76 S77 A78 A79 C80 K81 I82 S83 G84 C85 T86 F87 S88 A89 N90

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *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
X-PLOR	structure solution	3.851
X-PLOR	refinement	3.851

No chemical shift data was provided.

6 Model quality (i)

6.1 Standard geometry (i)

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

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	A	0.0±0.0	1.9±0.3
All	All	0	38

There are no bond-length outliers.

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	A	41	ARG	Sidechain	19
1	A	56	ARG	Sidechain	19

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	A	523	499	499	74±10
All	All	10460	9980	9980	1470

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 72.

5 of 531 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:45:VAL:HG21	1:A:57:ILE:HG22	1.03	1.26	14	10
1:A:27:VAL:HG13	1:A:42:PHE:CB	0.98	1.88	13	20
1:A:35:THR:HG23	1:A:50:SER:O	0.97	1.59	4	9
1:A:45:VAL:HG21	1:A:57:ILE:CG2	0.97	1.88	3	9
1:A:27:VAL:HG13	1:A:42:PHE:HB3	0.96	1.34	1	20

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	73/90 (81%)	44±2 (61±3%)	23±3 (31±4%)	6±2 (8±2%)	2 13
All	All	1460/1800 (81%)	889 (61%)	451 (31%)	120 (8%)	2 13

5 of 21 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	49	THR	17
1	A	66	THR	15
1	A	8	ASN	9
1	A	43	ASP	9
1	A	89	ALA	8

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	63/75 (84%)	43±3 (68±5%)	20±3 (32±5%)	1 14
All	All	1260/1500 (84%)	863 (68%)	397 (32%)	1 14

5 of 57 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	86	THR	17
1	A	58	SER	16
1	A	41	ARG	15
1	A	81	LYS	14
1	A	63	LYS	13

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