



Full wwPDB NMR Structure Validation Report ⓘ

Feb 20, 2018 – 03:49 pm GMT

PDB ID : 2K3P
Title : Solution structure of the C-terminal domain (TUSP1-C) of the egg case silk from *Nephila antipodiana*
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Deposited on : 2008-05-15

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 ⓘ symbol.

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	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	trunk30686
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30686

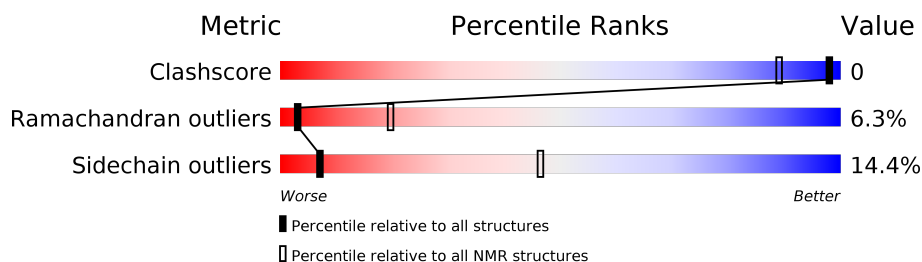
1 Overall quality at a glance

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	136279	12091
Ramachandran outliers	132675	10835
Sidechain outliers	132484	10811

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	161	

2 Ensemble composition and analysis

This entry contains 10 models. Model 9 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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:21 (19)	1.75	1
2	A:44-A:161 (118)	1.22	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 2 clusters and 2 single-model clusters were found.

Cluster number	Models
1	1, 4, 5, 6, 8
2	2, 9, 10
Single-model clusters	3; 7

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 2264 atoms, of which 1150 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called TuSp1.

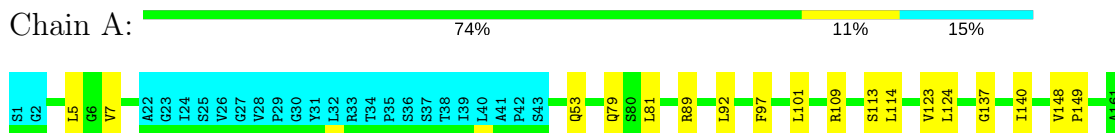
Mol	Chain	Residues	Atoms					Trace
1	A	161	Total	C	H	N	O	0
			2264	692	1150	190	232	

4 Residue-property plots

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: TuSp1

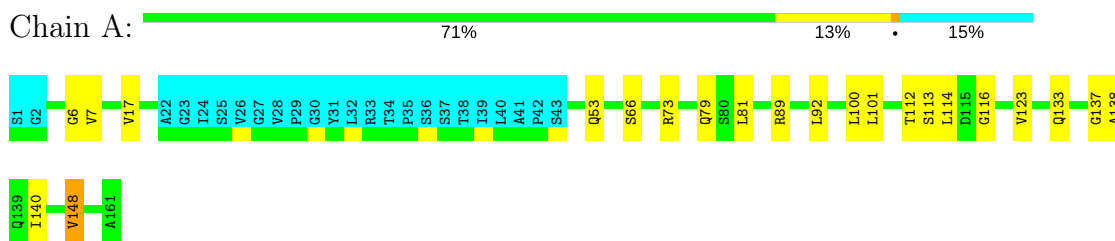


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

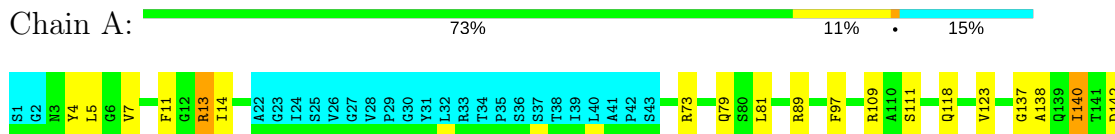
4.2.1 Score per residue for model 1

- Molecule 1: TuSp1



4.2.2 Score per residue for model 2

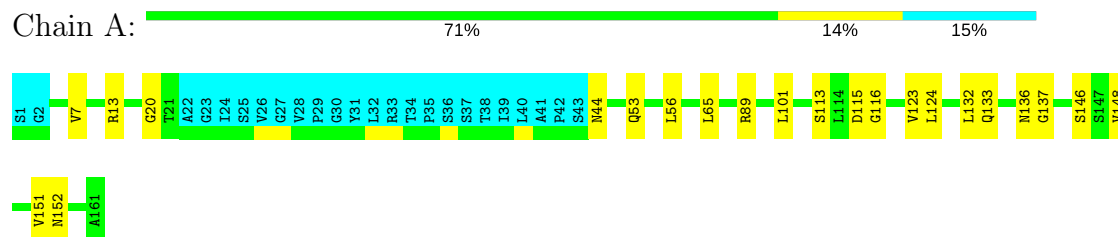
- Molecule 1: TuSp1





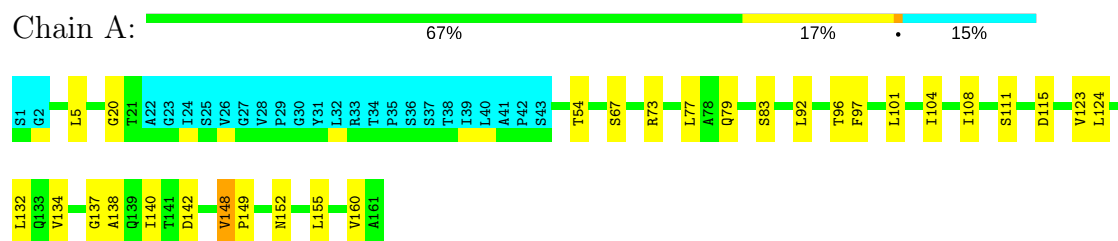
4.2.3 Score per residue for model 3

- Molecule 1: TuSp1



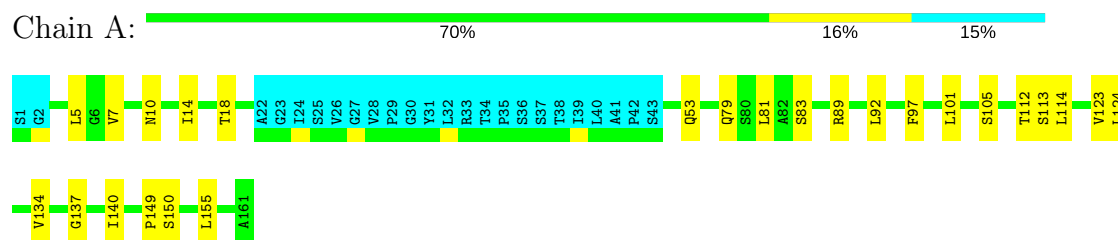
4.2.4 Score per residue for model 4

- Molecule 1: TuSp1



4.2.5 Score per residue for model 5

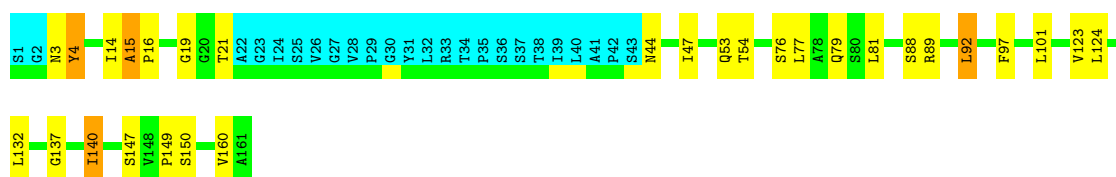
- Molecule 1: TuSp1



4.2.6 Score per residue for model 6

- Molecule 1: TuSp1

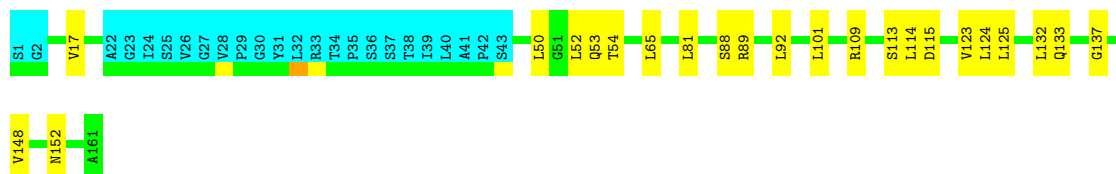




4.2.7 Score per residue for model 7

- Molecule 1: TuSp1

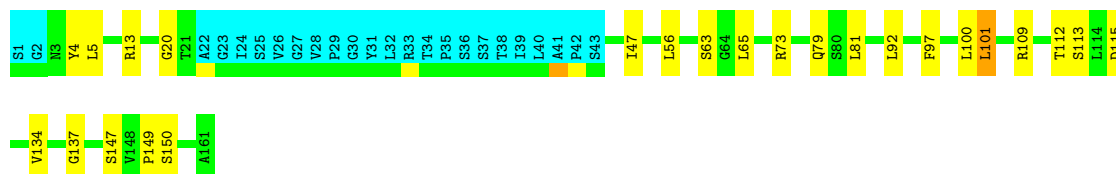
Chain A: 71% 14% 15%



4.2.8 Score per residue for model 8

- Molecule 1: TuSp1

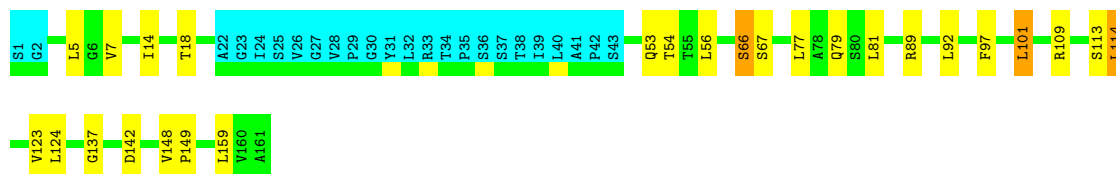
Chain A: 70% 14% 15%



4.2.9 Score per residue for model 9 (medoid)

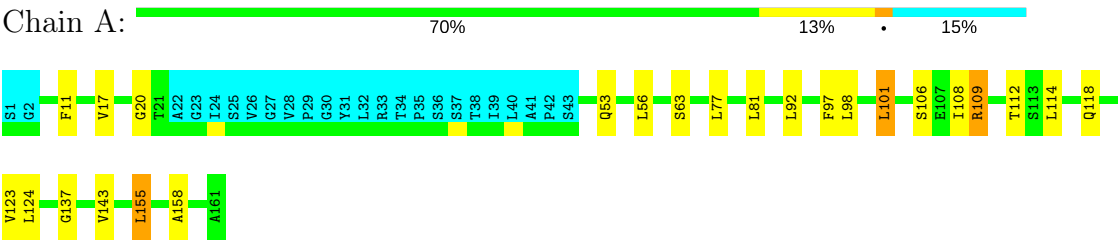
- Molecule 1: TuSp1

Chain A: 69% 14% 15%



4.2.10 Score per residue for model 10

- Molecule 1: TuSp1



5 Refinement protocol and experimental data overview

The models were refined using the following method: *torsion angle dynamics*.

Of the 100 calculated structures, 10 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
AMBER	refinement	7

No chemical shift data was provided. 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](#)

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	0.64±0.01	0±0/959 (0.0±0.0%)	1.04±0.03	0±0/1311 (0.0±0.0%)
All	All	0.64	0/9590 (0.0%)	1.04	3/13110 (0.0%)

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	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	73	ARG	NE-CZ-NH1	8.97	124.78	120.30	1	2
1	A	13	ARG	NE-CZ-NH1	5.03	122.82	120.30	2	1

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	A	954	987	994	1±1
All	All	9540	9870	9940	7

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 0.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:97:PHE:O	1:A:101:LEU:HD22	0.57	1.99	9	3
1:A:106:SER:HA	1:A:112:THR:HG21	0.54	1.79	10	1
1:A:15:ALA:HB1	1:A:16:PRO:HD2	0.51	1.82	6	1
1:A:15:ALA:HB1	1:A:16:PRO:CD	0.42	2.45	6	1
1:A:155:LEU:HA	1:A:158:ALA:HB3	0.40	1.94	10	1

6.3 Torsion angles

6.3.1 Protein backbone

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	136/161 (84%)	101±3 (74±2%)	26±4 (19±3%)	9±3 (6±2%)	3	20
All	All	1360/1610 (84%)	1010 (74%)	264 (19%)	86 (6%)	3	20

All 34 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	137	GLY	10
1	A	113	SER	6
1	A	140	ILE	5
1	A	7	VAL	5
1	A	149	PRO	5
1	A	14	ILE	4
1	A	148	VAL	4
1	A	20	GLY	4
1	A	17	VAL	3
1	A	138	ALA	3
1	A	92	LEU	3
1	A	67	SER	2
1	A	109	ARG	2
1	A	108	ILE	2
1	A	116	GLY	2
1	A	111	SER	2
1	A	18	THR	2
1	A	115	ASP	2

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Mol	Chain	Res	Type	Models (Total)
1	A	66	SER	2
1	A	4	TYR	2
1	A	89	ARG	2
1	A	147	SER	2
1	A	21	THR	1
1	A	3	ASN	1
1	A	15	ALA	1
1	A	63	SER	1
1	A	5	LEU	1
1	A	114	LEU	1
1	A	19	GLY	1
1	A	6	GLY	1
1	A	44	ASN	1
1	A	112	THR	1
1	A	143	VAL	1
1	A	150	SER	1

6.3.2 Protein sidechains ⓘ

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	109/127 (86%)	93±3 (86±2%)	16±3 (14±2%)	7	46
All	All	1090/1270 (86%)	933 (86%)	157 (14%)	7	46

All 53 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	123	VAL	9
1	A	101	LEU	9
1	A	81	LEU	8
1	A	53	GLN	7
1	A	124	LEU	7
1	A	79	GLN	7
1	A	92	LEU	6
1	A	89	ARG	5
1	A	114	LEU	5

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Mol	Chain	Res	Type	Models (Total)
1	A	132	LEU	4
1	A	5	LEU	4
1	A	97	PHE	4
1	A	77	LEU	4
1	A	109	ARG	4
1	A	54	THR	4
1	A	56	LEU	4
1	A	155	LEU	3
1	A	133	GLN	3
1	A	142	ASP	3
1	A	148	VAL	3
1	A	65	LEU	3
1	A	134	VAL	3
1	A	13	ARG	3
1	A	152	ASN	3
1	A	160	VAL	2
1	A	88	SER	2
1	A	140	ILE	2
1	A	115	ASP	2
1	A	47	ILE	2
1	A	100	LEU	2
1	A	4	TYR	2
1	A	83	SER	2
1	A	118	GLN	2
1	A	73	ARG	2
1	A	11	PHE	2
1	A	112	THR	2
1	A	150	SER	2
1	A	125	LEU	1
1	A	98	LEU	1
1	A	105	SER	1
1	A	146	SER	1
1	A	63	SER	1
1	A	136	ASN	1
1	A	151	VAL	1
1	A	159	LEU	1
1	A	50	LEU	1
1	A	66	SER	1
1	A	52	LEU	1
1	A	44	ASN	1
1	A	76	SER	1
1	A	96	THR	1

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Mol	Chain	Res	Type	Models (Total)
1	A	10	ASN	1
1	A	104	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](#)

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

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

No chemical shift data were provided