wwPDB NMR Structure Validation Summary Report (i)

Apr 26, 2016 – 08:29 PM BST

PDB ID	:	2G35
Title	:	NMR structure of talin-PTB in complex with PIPKI
Authors	:	Kong, X.; Wang, X.; Misra, S.; Qin, J.
Deposited on	:	2006-02-17

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:

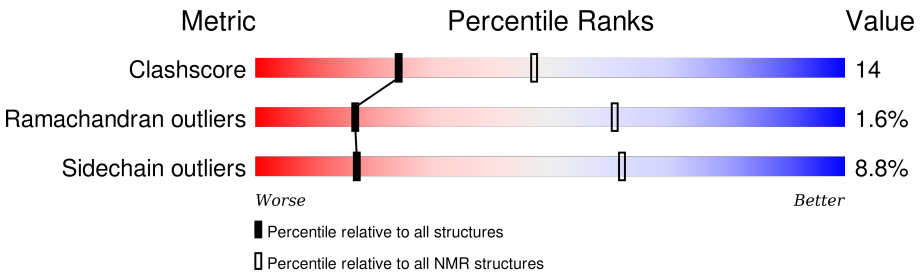
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	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 87%.

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	100	64% 19% • 15%
2	B	8	13% 50% 38%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 16 is the overall representative, medoid model (most similar to other models). The authors have identified model 11 as representative.

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:6-A:14, A:21-A:96, B:2-B:3, B:5-B:7 (90)	0.25	16

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 3 clusters and 5 single-model clusters were found.

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

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1805 atoms, of which 917 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Talin-1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	100	1666	528	853	133	149	3	0

- Molecule 2 is a protein called peptide.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
2	B	8	139	48	64	11	15	1	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: Talin-1

Chain A: 64% 19% • 15%

L1 K2 T3 Y4 G5 E13 K14 M15 K16 G17 K18 N19 K20 R24 T29 G32 V33 N34 R35 V36 E42 Q45 E46 W47 S58 P59 K60 L64 D65 D68 Y69 Q70 Y73 E80 G81 E82 Q83 K97 K98 K99 K100

- Molecule 2: peptide

Chain B: 13% 50% 38%

S1 W2 V3 Y4 S5 P6 L7 H8

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

The representative model is number 16. Colouring as in section 4.1 above.

- Molecule 1: Talin-1

Chain A: 51% 32% • 15%

L1 K2 T3 Y4 G5 F8 F9 L10 V11 K12 M15 K16 G17 K18 N19 K20 L25 T29 G32 V33 N34 R35 V36 D37 E38 K41 E42 Q45 E46 W47 R54 W55 A56 A57 S58 P59 K60 S61 F62 T63 L64 Q70 D71 G72 Y73 Q77 T78 T79 E80 G81

E82 Q83 I84 A85 Q86 L87 I88 Y91 K97 K98 K99 K100

- Molecule 2: peptide

Chain B: 63% 38%

S1 W2 V3 Y4 S5 P6 L7 H8

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *Distance geometry simulated annealing.*

Of the 50 calculated structures, 20 were deposited, based on the following criterion: *?*.

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

Software name	Classification	Version
X-PLOR	refinement	

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 7061
Number of chemical shift lists	2
Total number of shifts	1434
Number of shifts mapped to atoms	1371
Number of unparsed shifts	0
Number of shifts with mapping errors	63
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	87%

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: PTR

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	A	689	699	699	21±6
2	B	42	42	42	1±1
All	All	14620	14820	14820	416

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 14.

5 of 131 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:28:ILE:HD12	1:A:28:ILE:N	0.67	2.05	20	2
1:A:54:ARG:NH1	2:B:2:TRP:CH2	0.64	2.65	16	1
1:A:28:ILE:N	1:A:28:ILE:HD12	0.64	2.07	3	2
1:A:51:ASN:HD22	1:A:52:ILE:N	0.60	1.92	17	1
1:A:48:SER:N	1:A:51:ASN:OD1	0.59	2.35	17	1

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	85/100 (85%)	80±1 (94±1%)	5±1 (6±2%)	0±0 (0±1%)	38	79
2	B	5/8 (62%)	4±1 (70±13%)	1±1 (10±13%)	1±1 (20±15%)	0	2
All	All	1800/2160 (83%)	1660 (92%)	112 (6%)	28 (2%)	17	61

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)
2	B	7	LEU	11
1	A	59	PRO	7
2	B	2	TRP	6
2	B	6	PRO	3
1	A	81	GLY	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	A	76/89 (85%)	69±2 (91±3%)	7±2 (9±3%)	18	63
2	B	5/7 (71%)	4±0 (88±10%)	1±0 (12±10%)	11	53
All	All	1620/1920 (84%)	1477 (91%)	143 (9%)	17	62

5 of 26 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	32	CYS	20
1	A	58	SER	14

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Mol	Chain	Res	Type	Models (Total)
1	A	42	GLU	14
1	A	64	LEU	13
2	B	5	SER	10

6.3.3 RNA (i)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains (i)

1 non-standard protein/DNA/RNA residue is 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
2	PTR	B	4	2	13,16,17	1.33±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.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
2	PTR	B	4	2	19,22,24	1.44±0.02	1±0 (5±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
2	PTR	B	4	2	-	0±0,9,11,13	0±0,1,1,1

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
2	B	4	PTR	P-OH-CZ	5.26	108.92	123.85	4	20

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 87% for the well-defined parts and 87% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 7061

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	1333
Number of shifts mapped to atoms	1270
Number of unparsed shifts	0
Number of shifts with mapping errors	63
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	6

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

- Residue not found in structure. First 5 (of 63) occurences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
A	102	LYS	HE2	2.98	-1.0	1
A	102	LYS	HA	4.3	-1.0	1
A	105	PHE	HB2	2.98	-1.0	1
A	103	ASP	H	8.15	-1.0	1
A	105	PHE	HE1	7.17	-1.0	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
$^{13}C_{\alpha}$	106	-0.22 ± 0.12	None needed (< 0.5 ppm)
$^{13}C_{\beta}$	98	-0.05 ± 0.16	None needed (< 0.5 ppm)
$^{13}C'$	0	—	—

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Nucleus	# values	Correction ± precision, *ppm*	Suggested action
^{15}N	103	-0.29 ± 0.62	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 82%, i.e. 933 atoms were assigned a chemical shift out of a possible 1141. 15 out of 17 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	336/444 (76%)	168/177 (95%)	85/180 (47%)	83/87 (95%)
Sidechain	505/593 (85%)	309/345 (90%)	187/226 (83%)	9/22 (41%)
Aromatic	92/104 (88%)	48/54 (89%)	42/47 (89%)	2/3 (67%)
Overall	933/1141 (82%)	525/576 (91%)	314/453 (69%)	94/112 (84%)

7.1.4 Statistically unusual chemical shifts (i)

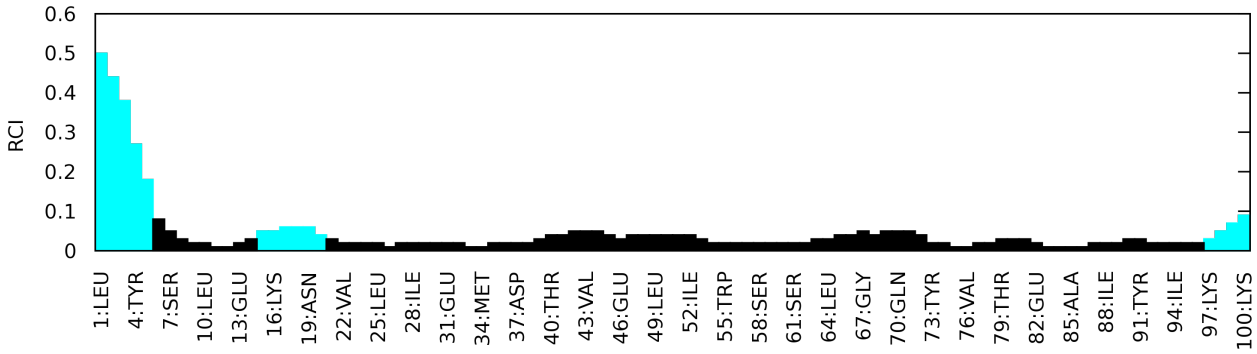
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
1	A	30	LYS	HD2	0.03	2.76 – 0.46	-6.9
1	A	41	LYS	HD2	0.03	2.76 – 0.46	-6.9
1	A	14	LYS	HD2	0.03	2.76 – 0.46	-6.9
1	A	84	ILE	HG12	-1.24	3.27 – -0.73	-6.3
1	A	45	GLN	HG3	0.77	3.75 – 0.85	-5.3
1	A	1	LEU	HB2	-0.10	3.32 – -0.08	-5.1

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:



7.2 Chemical shift list 2

File name: BMRB entry 7061

Chemical shift list name: *assigned_chem_shift_list_2*

7.2.1 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	101
Number of shifts mapped to atoms	101
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

7.2.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	8	0.69 ± 0.74	None needed (imprecise)
$^{13}C_{\beta}$	8	0.00 ± 0.00	None needed (< 0.5 ppm)
$^{13}C'$	0	—	—
^{15}N	7	-0.03 ± 0.48	None needed (< 0.5 ppm)

7.2.3 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 5%, i.e. 60 atoms were assigned a chemical shift out of a possible 1141. 2 out of 17 assigned methyl groups (LEU and VAL) were assigned

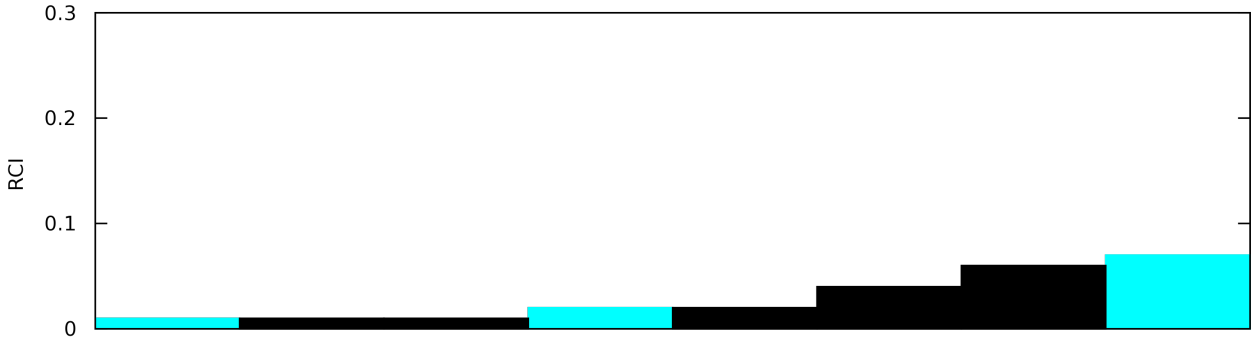
stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	18/444 (4%)	9/177 (5%)	5/180 (3%)	4/87 (5%)
Sidechain	30/593 (5%)	18/345 (5%)	12/226 (5%)	0/22 (0%)
Aromatic	12/104 (12%)	6/54 (11%)	5/47 (11%)	1/3 (33%)
Overall	60/1141 (5%)	33/576 (6%)	22/453 (5%)	5/112 (4%)

7.2.4 Statistically unusual chemical shifts (i)

There are no statistically unusual chemical shifts.

7.2.5 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 B: