wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1XKE
Title	:	Solution structure of the second Ran-binding domain from human RanBP2
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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
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:

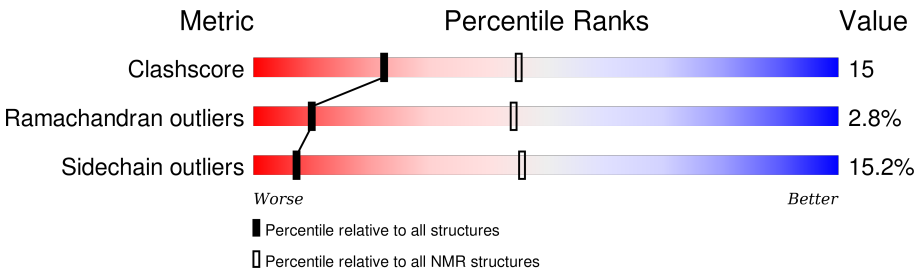
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 69%.

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	130	48% 36% 7% 9%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 2 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:5-A:122 (118)	0.55	2

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. No single-model clusters were found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2142 atoms, of which 1080 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Ran-binding protein 2.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	130	2142	674	1080	186	196	6	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	CLONING ARTIFACT	UNP P49792
A	2	SER	-	CLONING ARTIFACT	UNP P49792
A	130	LYS	-	CLONING ARTIFACT	UNP P49792

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: Ran-binding protein 2

Chain A: 48% 36% 7% 9%

G1 S2 G3 E4 E5 D6 E7 K8 S12 R19 F20 D21 V24 E29 R30 G33 N34 L35 K36 I37 L38 L46 R47 M48 L49 M50 R51 R52 E53 Q54 V55 L56 K57 V58 C59 A60 N61 H62 W63 I64 T65 T66 T67 M68 N69 L70 L73 D77 R78 A79 W80 M81

W82 L83 A84 S85 D86 F87 G90 D91 A92 K93 L94 E95 Q96 L97 K100 P104 A107 E108 K111 Q118 R119 L120 D123 I124 P125 L126 Q127 T128 P129 K130

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

The representative model is number 2. Colouring as in section 4.1 above.

- Molecule 1: Ran-binding protein 2

Chain A: 46% 31% 12% • 9%

G1 S2 G3 E4 E5 D6 E7 S12 F18 D21 V24 E29 G33 N34 L35 L46 R47 M48 L49 M50 R51 R52 E53 Q54 V55 L56 K57 V58 C59 A60 N61 I64 T65 T66 T67 M68 N69 L70 L73 D77 R78 A79 W80 M81 W82 L83 A84 S85 D86 F87

D91 A92 K93 L94 E95 Q96 L97 A98 A99 K100 F101 P104 A107 E108 E109 F110 K111 Q112 K113 F114 E115 Q118 R119 L120 L121 L122 D123 I124 P125 L126 Q127 T128 P129 K130

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *Simulated annealing.*

Of the 200 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
CNS	structure solution	1.1
CNS	refinement	1.1
XPLOR-NIH	structure solution	2.9.6

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 5159
Number of chemical shift lists	1
Total number of shifts	1308
Number of shifts mapped to atoms	1308
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	69%

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	1.50±0.03	13±3/993 (1.3±0.3%)	1.03±0.02	0±0/1331 (0.0±0.0%)
All	All	1.50	252/19860 (1.3%)	1.04	5/26620 (0.0%)

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	0.7±0.8
All	All	0	14

5 of 36 unique bond 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(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	50	MET	N-CA	-7.88	1.30	1.46	13	20
1	A	50	MET	C-N	-7.10	1.17	1.34	14	20
1	A	96	GLN	N-CA	-7.08	1.32	1.46	20	20
1	A	5	GLU	N-CA	-6.73	1.32	1.46	17	4
1	A	90	GLY	N-CA	-6.46	1.36	1.46	13	4

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})	Models	
								Worst	Total
1	A	56	LEU	N-CA-CB	-5.38	99.65	110.40	10	5

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	51	ARG	Sidechain	6
1	A	30	ARG	Sidechain	4
1	A	78	ARG	Sidechain	2
1	A	19	ARG	Sidechain	1
1	A	119	ARG	Sidechain	1

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	975	993	991	29±3
All	All	19500	19860	19817	589

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 15.

5 of 205 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:17:LEU:HB2	1:A:50:MET:SD	0.81	2.15	10	4
1:A:76:SER:HB3	1:A:79:ALA:HB3	0.75	1.56	13	2
1:A:60:ALA:HA	1:A:86:ASP:OD1	0.71	1.86	7	10
1:A:18:PHE:O	1:A:97:LEU:HB2	0.69	1.88	13	3
1:A:19:ARG:HA	1:A:97:LEU:HD22	0.68	1.65	11	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	118/130 (91%)	109±1 (92±1%)	6±2 (5±1%)	3±2 (3±1%)	10	44
All	All	2360/2600 (91%)	2181 (92%)	114 (5%)	65 (3%)	10	44

5 of 11 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	58	VAL	14
1	A	87	PHE	12
1	A	78	ARG	7
1	A	53	GLU	7
1	A	89	ASP	7

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	105/115 (91%)	89±3 (85±3%)	16±3 (15±3%)	7 46
All	All	2100/2300 (91%)	1781 (85%)	319 (15%)	7 46

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	85	SER	19
1	A	64	ILE	19
1	A	5	GLU	17
1	A	119	ARG	16
1	A	46	LEU	16

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 (i)

The completeness of assignment taking into account all chemical shift lists is 69% for the well-defined parts and 71% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 5159

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	1308
Number of shifts mapped to atoms	1308
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	2

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	129	-0.32 ± 0.10	None needed (< 0.5 ppm)
$^{13}C_{\beta}$	123	-0.07 ± 0.15	None needed (< 0.5 ppm)
$^{13}C'$	122	-0.08 ± 0.16	None needed (< 0.5 ppm)
^{15}N	124	0.79 ± 0.34	Should be applied

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 69%, i.e. 1089 atoms were assigned a chemical shift out of a possible 1568. 1 out of 23 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	582/586 (99%)	234/234 (100%)	232/236 (98%)	116/116 (100%)
Sidechain	507/865 (59%)	326/509 (64%)	172/308 (56%)	9/48 (19%)

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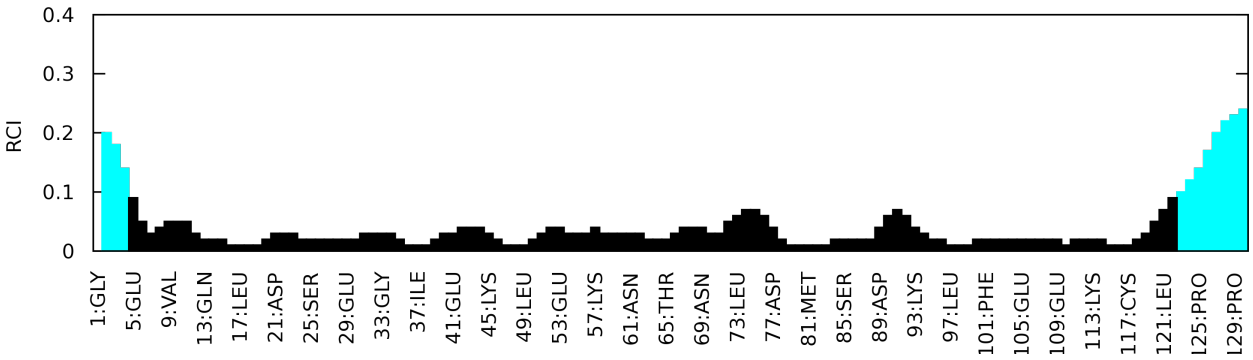
	Total	^{1}H	^{13}C	^{15}N
Aromatic	0/117 (0%)	0/62 (0%)	0/50 (0%)	0/5 (0%)
Overall	1089/1568 (69%)	560/805 (70%)	404/594 (68%)	125/169 (74%)

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	119	ARG	NE	123.40	92.63 – 76.73	24.4
1	A	20	PHE	HB3	0.99	4.85 – 1.05	-5.2

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: