wwPDB NMR Structure Validation Summary Report (i)

Apr 26, 2016 – 07:54 PM BST

PDB ID	:	2EHR
Title	:	Solution structure of the sixth PDZ domain of human InaD-like protein
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Deposited on	:	2007-03-07

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:

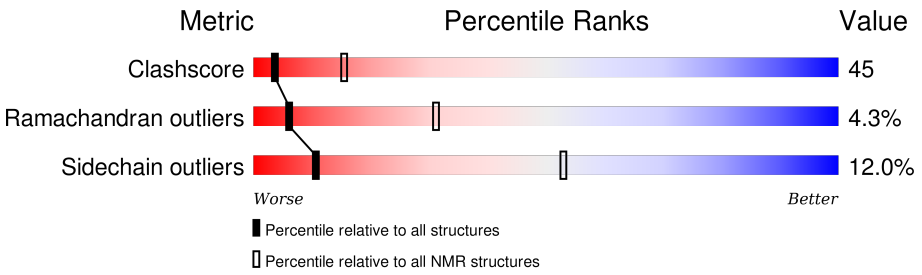
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 95%.

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	117	25% 39% • 32%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 3 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:14-A:35, A:52-A:108 (79)	0.13	3

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1758 atoms, of which 890 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called InaD-like protein.

Mol	Chain	Residues	Atoms					Trace
			Total	C	H	N	O	
1	A	117	1758	547	890	153	168	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	EXPRESSION TAG	UNP Q8NI35
A	2	SER	-	EXPRESSION TAG	UNP Q8NI35
A	3	SER	-	EXPRESSION TAG	UNP Q8NI35
A	4	GLY	-	EXPRESSION TAG	UNP Q8NI35
A	5	SER	-	EXPRESSION TAG	UNP Q8NI35
A	6	SER	-	EXPRESSION TAG	UNP Q8NI35
A	7	GLY	-	EXPRESSION TAG	UNP Q8NI35

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: InaD-like protein

Chain A: 25% 39% • 32%

G1 S2 S3 G4 S5 S6 G7 P8 N9 F10 S11 H12 W13 I18 I21 F22 R23 E24 P25 N26 V27 S28 L29 G30 I31 S32 I33 V34 G35 G36 Q37 T38 V39 I40 K41 R42 L43 K44 N45 G46 E47 E48 L49 K50 G51 I52 F53 I54 K55 Q56 V57 L58 S61 P62 A63 T66

N67 A68 L69 K70 D73 K74 I75 L76 E77 V78 S79 G80 V81 D82 L83 A86 S87 H88 A91 V92 E93 A94 I95 K96 N97 A98 G99 N100 P101 V102 V103 F104 I105 V106 Q107 S108 L109 S110 S111 T112 P113 R114 V115 I116 P117

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

The representative model is number 3. Colouring as in section 4.1 above.

- Molecule 1: InaD-like protein

Chain A: 23% 40% • 32%

G1 S2 S3 G4 S5 S6 G7 P8 N9 F10 S11 H12 W13 G14 I18 I21 F22 R23 E24 P25 N26 V27 S28 L29 G30 I31 S32 I33 V34 G35 G36 Q37 T38 V39 I40 K41 R42 L43 K44 N45 G46 E47 E48 L49 K50 G51 I52 F53 I54 K55 Q56 V57 L58 S61 P62 A63

T66 N67 A68 L69 K70 D73 K74 I75 L76 E77 V78 S79 G80 V81 D82 L83 A86 S87 H88 A91 V92 E93 A94 I95 K96 N97 A98 G99 N100 P101 V102 V103 F104 I105 V106 Q107 S108 L109 S110 S111 T112 P113 R114 V115 I116 P117

5 Refinement protocol and experimental data overview (i)

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

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *structures with the least restraint violations, structures with the lowest energy, target function.*

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

Software name	Classification	Version
CYANA	structure solution	2.1
CYANA	refinement	2.1

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 11197
Number of chemical shift lists	1
Total number of shifts	1384
Number of shifts mapped to atoms	1384
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	95%

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)

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	586	606	606	53±6
All	All	11720	12120	12120	1067

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 45.

5 of 159 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:33:ILE:HD12	1:A:88:HIS:CD2	0.97	1.94	4	6
1:A:21:ILE:HD12	1:A:104:PHE:CD1	0.96	1.95	16	18
1:A:17:ARG:NH1	1:A:106:VAL:HG23	0.91	1.80	8	1
1:A:21:ILE:HD11	1:A:69:LEU:HD11	0.90	1.43	19	14
1:A:52:ILE:HD11	1:A:86:ALA:O	0.89	1.67	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	79/117 (68%)	67±1 (84±2%)	9±1 (11±2%)	3±1 (4±1%)	6 31
All	All	1580/2340 (68%)	1331 (84%)	181 (11%)	68 (4%)	6 31

5 of 7 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	99	GLY	20
1	A	25	PRO	20
1	A	67	ASN	15
1	A	73	ASP	5
1	A	35	GLY	4

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	66/98 (67%)	58±2 (88±2%)	8±2 (12±2%)	11 53
All	All	1320/1960 (67%)	1162 (88%)	158 (12%)	11 53

5 of 21 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	34	VAL	20
1	A	61	SER	20
1	A	78	VAL	20
1	A	76	LEU	20
1	A	56	GLN	15

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 95% for the well-defined parts and 89% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 11197

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	1384
Number of shifts mapped to atoms	1384
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

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	108	-0.06 ± 0.10	None needed (< 0.5 ppm)
$^{13}C_{\beta}$	99	-0.08 ± 0.13	None needed (< 0.5 ppm)
$^{13}C'$	105	-0.10 ± 0.13	None needed (< 0.5 ppm)
^{15}N	97	0.05 ± 0.56	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 95%, i.e. 877 atoms were assigned a chemical shift out of a possible 923. 0 out of 15 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	383/385 (99%)	153/153 (100%)	156/158 (99%)	74/74 (100%)
Sidechain	464/504 (92%)	282/292 (97%)	172/193 (89%)	10/19 (53%)

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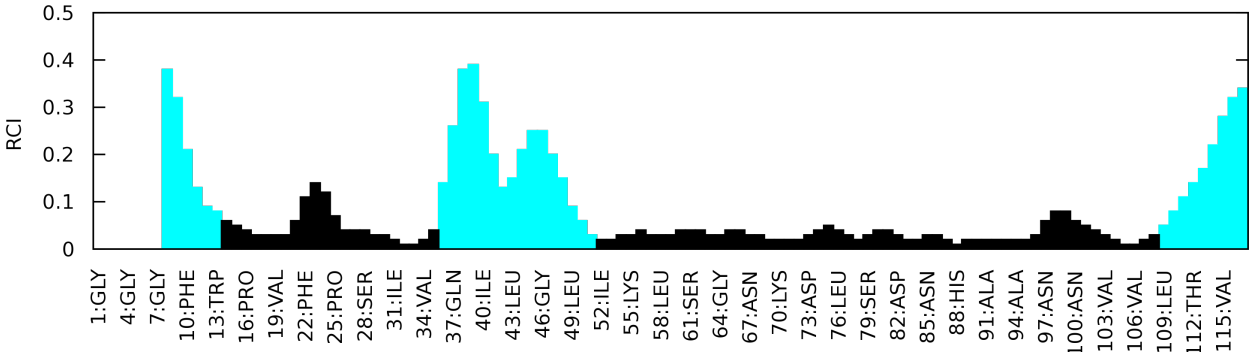
	Total	^{1}H	^{13}C	^{15}N
Aromatic	30/34 (88%)	17/19 (89%)	13/14 (93%)	0/1 (0%)
Overall	877/923 (95%)	452/464 (97%)	341/365 (93%)	84/94 (89%)

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: