Full wwPDB NMR Structure Validation Report (i)

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PDB ID	:	2RNA
Title	:	Itk SH3 average minimized
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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
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:

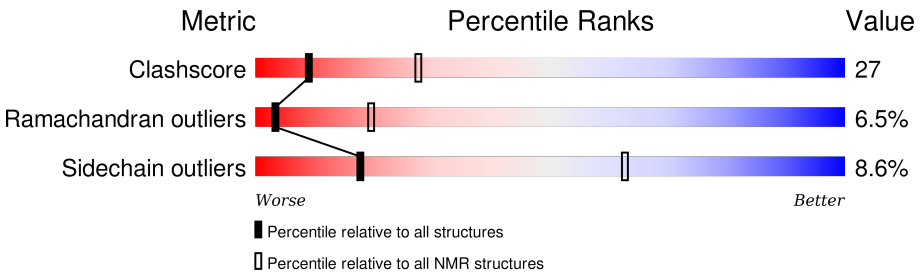
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 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	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	64	63% 31% 6%

2 Ensemble composition and analysis (i)

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1007 atoms, of which 479 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Tyrosine-protein kinase ITK/TSK.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	64	1007	332	479	84	111	1	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	GLY	-	EXPRESSION TAG	UNP Q03526
A	170	SER	-	EXPRESSION TAG	UNP Q03526

4 Residue-property plots (i)

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: Tyrosine-protein kinase ITK/TSK

Chain A:

5 Refinement protocol and experimental data overview (i)

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

Of the ? calculated structures, 1 were deposited, based on the following criterion: *?*.

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

Software name	Classification	Version
CNS	refinement	

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)

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	528	479	476	27
All	All	528	479	476	27

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 27.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:171:PRO:O	1:A:173:GLU:N	0.80	2.13
1:A:205:GLU:CG	1:A:206:ILE:N	0.72	2.53
1:A:184:THR:HG23	1:A:186:ASP:H	0.62	1.53
1:A:205:GLU:OE1	1:A:208:TRP:C	0.59	2.40
1:A:205:GLU:OE1	1:A:208:TRP:O	0.56	2.22
1:A:205:GLU:HG2	1:A:206:ILE:N	0.56	2.16
1:A:205:GLU:CG	1:A:206:ILE:H	0.56	2.13
1:A:183:GLN:NE2	1:A:184:THR:O	0.55	2.40
1:A:205:GLU:HG3	1:A:206:ILE:H	0.52	1.64
1:A:205:GLU:CD	1:A:208:TRP:C	0.50	2.70
1:A:204:SER:O	1:A:205:GLU:O	0.50	2.29
1:A:205:GLU:OE2	1:A:208:TRP:C	0.49	2.51
1:A:214:LYS:CD	1:A:214:LYS:N	0.48	2.77
1:A:205:GLU:OE2	1:A:208:TRP:CA	0.48	2.62
1:A:183:GLN:NE2	1:A:184:THR:H	0.48	2.07

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:183:GLN:NE2	1:A:184:THR:N	0.47	2.63
1:A:229:LYS:O	1:A:230:SER:O	0.47	2.33
1:A:171:PRO:O	1:A:173:GLU:O	0.45	2.34
1:A:183:GLN:HE21	1:A:184:THR:N	0.45	2.09
1:A:214:LYS:CD	1:A:214:LYS:H	0.43	2.26
1:A:194:CYS:O	1:A:195:ASP:CB	0.43	2.66
1:A:205:GLU:OE2	1:A:208:TRP:N	0.43	2.52
1:A:205:GLU:OE2	1:A:209:TRP:N	0.42	2.52
1:A:184:THR:HG23	1:A:185:ASN:N	0.41	2.30
1:A:185:ASN:C	1:A:186:ASP:OD1	0.41	2.58
1:A:188:GLN:O	1:A:220:TYR:N	0.40	2.52
1:A:171:PRO:O	1:A:172:GLU:C	0.40	2.57

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	62/64 (97%)	54 (87%)	4 (6%)	4 (6%)	3 19
All	All	62/64 (97%)	54 (87%)	4 (6%)	4 (6%)	3 19

All 4 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	172	GLU
1	A	170	SER
1	A	230	SER
1	A	205	GLU

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	58/58 (100%)	53 (91%)	5 (9%)	18	63
All	All	58/58 (100%)	53 (91%)	5 (9%)	18	63

All 5 residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	184	THR
1	A	199	TYR
1	A	227	VAL
1	A	193	ARG
1	A	189	GLU

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)

No chemical shift data were provided