wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	2I3B
Title	:	Solution Structure of a Human Cancer-Related Nucleoside Triphosphatase
Authors	:	Placzek, W.J.; Almeida, M.S.; Wuthrich, K.
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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:

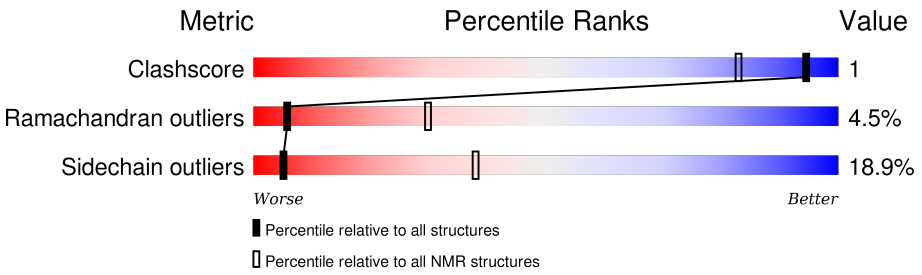
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 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	189	79% 15% • 6%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 1 is the overall representative, medoid model (most similar to other models).

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:2-A:97, A:106-A:187 (178)	0.72	1

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 5 clusters and 2 single-model clusters were found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2954 atoms, of which 1508 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Human Cancer-Related NTPase.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	189	2954	912	1508	264	265	5	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: Human Cancer-Related NTPase

Chain A: 79% 15% • 6%

A1 R2 V12 G13 K14 L17 R45 R46 V51 R58 L61 S62 R63 V64 K72 R73 R76 Q79 Y80 V81 V82 L90 A91 L92 R96 N97 A98 D99 C100 S101 S102 G103 P104 G105 E113 F124 K149 K161 K164 N167 V168 L176

S187 R188 K189

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

The representative model is number 1. Colouring as in section 4.1 above.

- Molecule 1: Human Cancer-Related NTPase

Chain A: 68% 23% • 6%

A1 R2 H3 V4 F5 G11 V12 L17 D33 Y36 T37 E38 R45 V51 T57 R58 G59 R63 L66 K72 R76 V77 G78 V82 D83 E88 Q89 L90 A91 L92 L95 R96 N97 A98 D99 C100 S101 S102 G103 P104 G105 E113 I114 G115

E118 F124 Q130 S133 V145 P146 K147 G148 K149 E155 E156 I157 K161 D162 V163 K164 V165 F166 N167 V168 L176 L177 P178 D179 Q185 S186 S187 R188 K189

5 Refinement protocol and experimental data overview (i)

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

Of the 80 calculated structures, 20 were deposited, based on the following criterion: *target function*.

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

Software name	Classification	Version
CYANA	structure solution	version 1.0.3
CYANA	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)

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	0.56±0.01	0±0/1400 (0.0±0.0%)	1.11±0.03	3±2/1899 (0.2±0.1%)
All	All	0.56	0/28000 (0.0%)	1.11	63/37980 (0.2%)

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	2.5±1.2
All	All	0	50

There are no bond-length outliers.

5 of 38 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(^{o})	Ideal(^{o})	Models	
								Worst	Total
1	A	63	ARG	NE-CZ-NH2	-7.25	116.67	120.30	15	4
1	A	51	VAL	CA-CB-CG2	7.03	121.45	110.90	3	4
1	A	168	VAL	CA-CB-CG2	6.96	121.35	110.90	5	1
1	A	2	ARG	NE-CZ-NH2	-6.79	116.90	120.30	20	3
1	A	134	THR	CA-CB-CG2	6.79	121.90	112.40	7	1

There are no chirality outliers.

5 of 23 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	76	ARG	Sidechain	6
1	A	80	TYR	Sidechain	5
1	A	45	ARG	Sidechain	4

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Mol	Chain	Res	Type	Group	Models (Total)
1	A	36	TYR	Sidechain	3
1	A	63	ARG	Sidechain	3

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	1375	1440	1440	4±2
All	All	27500	28800	28800	83

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 1.

5 of 59 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:6:LEU:HD21	1:A:180:ILE:HD11	0.81	1.51	9	2
1:A:6:LEU:HD21	1:A:18:ILE:HD11	0.71	1.62	17	1
1:A:6:LEU:HD11	1:A:180:ILE:HD11	0.69	1.64	4	4
1:A:92:LEU:HD21	1:A:123:LEU:HD11	0.68	1.64	7	1
1:A:64:VAL:HG12	1:A:65:GLY:H	0.62	1.55	12	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	178/189 (94%)	142±5 (80±3%)	28±5 (16±3%)	8±2 (4±1%)	6 30
All	All	3560/3780 (94%)	2837 (80%)	563 (16%)	160 (4%)	6 30

5 of 44 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	92	LEU	18
1	A	12	VAL	13
1	A	115	GLY	8
1	A	79	GLN	7
1	A	113	GLU	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	157/164 (96%)	127±5 (81±3%)	30±5 (19±3%)	5 38
All	All	3140/3280 (96%)	2545 (81%)	595 (19%)	5 38

5 of 107 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	149	LYS	20
1	A	90	LEU	19
1	A	92	LEU	18
1	A	167	ASN	18
1	A	124	PHE	17

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