# Full wwPDB NMR Structure Validation Report (i)

Apr 26, 2016 – 02:35 PM BST

| | | |
|---|---|---|
| PDB ID | : | 1FV5 |
| Title | : | SOLUTION STRUCTURE OF THE FIRST ZINC FINGER FROM THE DROSOPHILA U-SHAPED TRANSCRIPTION FACTOR |
| Authors | : | Liew, C.K.; Kowalski, K.; Fox, A.H.; Newton, A.; Sharpe, B.K.; Crossley, M.; Mackay, J.P. |
| Deposited on | : | 2000-09-18 |

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:

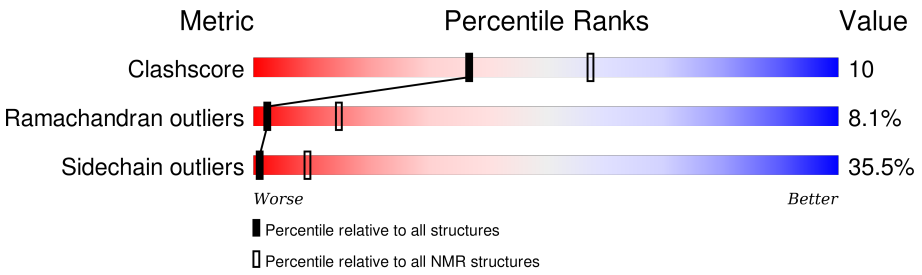
| | | |
|---|---|---|
| Cyrange | : | Kirchner and Güntert (2011) |
| NmrClust | : | Kelley et al. (1996) |
| MolProbity | : | 4.02b-467 |
| Mogul | : | unknown |
| Percentile statistics | : | 20151230.v01 (using entries in the PDB archive December 30th 2015) |
| RCI | : | v_1n_11_5_13_A (Berjanski et 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 is 50%.

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 | 36 | 50% 39% 8% • |

# 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:36 (35) | 0.38 | 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 3 clusters and 5 single-model clusters were found.

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

# 3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 549 atoms, of which 272 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called FIRST ZINC FINGER OF U-SHAPED.

| Mol | Chain | Residues | Atoms | | | | | | Trace |
|---|---|---|---|---|---|---|---|---|---|
| | | | Total | C | H | N | O | S | |
| 1 | A | 36 | 548 | 176 | 272 | 48 | 48 | 4 | 0 |

There are 2 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|---|---|---|---|---|---|
| A | 1 | GLY | - | cloning artifact | UNP Q9VPQ6 |
| A | 2 | SER | - | cloning artifact | UNP Q9VPQ6 |

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

| Mol | Chain | Residues | Atoms | |
|---|---|---|---|---|
| | | | Total | Zn |
| 2 | A | 1 | 1 | 1 |

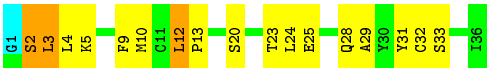
# 4 Residue-property plots (i)

## 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: FIRST ZINC FINGER OF U-SHAPED

Chain A:



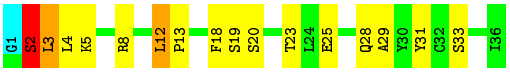

## 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

### 4.2.1 Score per residue for model 1 (medoid)

• Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A:




### 4.2.2 Score per residue for model 2

• Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A:

### 4.2.3 Score per residue for model 3

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 44% 39% 14% •

G1 S2 L3 L4 K5 P6 A7 R8 F9 M10 C11 L12 P13 C14 G15 I16 A17 S20 L24 E25 Q28 Y31 I36

### 4.2.4 Score per residue for model 4

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 61% 25% 11% •

G1 S2 L3 L4 F9 L12 P13 S20 P21 E25 Q28 A29 Y30 Y31 C32 S33 I36

### 4.2.5 Score per residue for model 5

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 44% 39% 11% • •

G1 S2 L3 L4 A7 R8 F9 M10 C11 L12 P13 A17 F18 S19 S20 P21 S22 T23 L24 H27 Q28 Y31 C32 S33 I36

### 4.2.6 Score per residue for model 6

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 44% 44% 8% •

G1 S2 L3 L4 K5 R8 F9 M10 C11 L12 P13 C14 A17 F18 S19 S20 P21 S22 T23 L24 Q28 C32 S33 I36

### 4.2.7 Score per residue for model 7

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

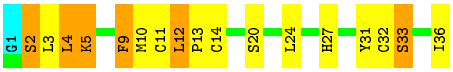
Chain A: 47% 42% 8% •

G1 S2 L3 L4 K5 P6 A7 M10 C11 L12 P13 S19 S20 P21 S22 T23 L24 E25 Q28 A29 Y30 Y31 C32 S33 I36

### 4.2.8 Score per residue for model 8

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 50% 31% 17% •




### 4.2.9 Score per residue for model 9

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 50% 36% 6% 6% •

### 4.2.10 Score per residue for model 10

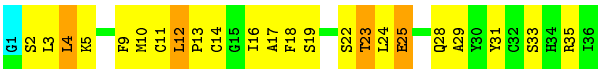
- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 47% 39% 8% • •

### 4.2.11 Score per residue for model 11

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 33% 53% 11% •




### 4.2.12 Score per residue for model 12

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 47% 36% 11% • •

### 4.2.13 Score per residue for model 13

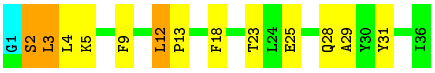
- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 47% 36% 11%

### 4.2.14 Score per residue for model 14

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

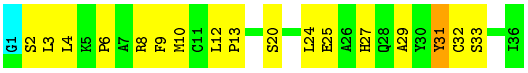
Chain A: 61% 28% 8%




### 4.2.15 Score per residue for model 15

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 50% 44%




### 4.2.16 Score per residue for model 16

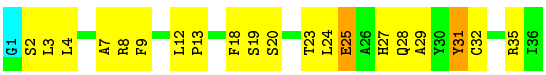
- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A: 44% 39% 14%

### 4.2.17 Score per residue for model 17

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

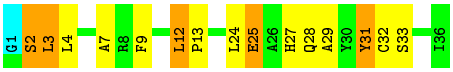
Chain A: 42% 50% 6%



### 4.2.18 Score per residue for model 18

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A:




### 4.2.19 Score per residue for model 19

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A:

### 4.2.20 Score per residue for model 20

- Molecule 1: FIRST ZINC FINGER OF U-SHAPED

Chain A:

# 5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *distance geometry simulated annealing cartesian dynamics.*

Of the 500 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 |
|---|---|---|
| DYANA | structure solution | 1.5 |
| CNS | refinement | 1.0 |

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 4644 |
|---|---|
| Number of chemical shift lists | 1 |
| Total number of shifts | 247 |
| Number of shifts mapped to atoms | 247 |
| Number of unparsed shifts | 0 |
| Number of shifts with mapping errors | 0 |
| Number of shifts with mapping warnings | 0 |
| Assignment completeness (well-defined parts) | 50% |

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: ZN

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 | 272 | 267 | 266 | 5±1 |
| All | All | 5460 | 5340 | 5320 | 108 |

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 10.

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

| Atom-1 | Atom-2 | Clash(Å) | Distance(Å) | Models | |
|---|---|---|---|---|---|
| | | | | Worst | Total |
| 1:A:9:PHE:HA | 1:A:24:LEU:HD11 | 0.67 | 1.66 | 6 | 3 |
| 1:A:9:PHE:CB | 1:A:24:LEU:HD11 | 0.63 | 2.22 | 11 | 4 |
| 1:A:2:SER:HB3 | 1:A:3:LEU:HD23 | 0.62 | 1.71 | 1 | 7 |
| 1:A:16:ILE:HD11 | 1:A:27:HIS:CE1 | 0.61 | 2.30 | 16 | 1 |
| 1:A:25:GLU:O | 1:A:29:ALA:HB2 | 0.59 | 1.97 | 17 | 12 |
| 1:A:10:MET:SD | 1:A:12:LEU:HD23 | 0.57 | 2.38 | 6 | 1 |
| 1:A:9:PHE:HB3 | 1:A:24:LEU:HD11 | 0.57 | 1.76 | 3 | 4 |
| 1:A:4:LEU:HD13 | 1:A:4:LEU:O | 0.55 | 2.02 | 9 | 1 |
| 1:A:12:LEU:N | 1:A:13:PRO:HD2 | 0.54 | 2.18 | 7 | 9 |
| 1:A:27:HIS:HA | 1:A:31:TYR:HB2 | 0.53 | 1.80 | 17 | 8 |
| 1:A:2:SER:C | 1:A:3:LEU:HD23 | 0.53 | 2.24 | 5 | 1 |
| 1:A:9:PHE:HB2 | 1:A:24:LEU:HD11 | 0.53 | 1.79 | 11 | 1 |

*Continued on next page...*

*Continued from previous page...*

| Atom-1 | Atom-2 | Clash(Å) | Distance(Å) | Models | |
|---|---|---|---|---|---|
| | | | | Worst | Total |
| 1:A:12:LEU:N | 1:A:13:PRO:CD | 0.50 | 2.74 | 1 | 20 |
| 1:A:20:SER:CB | 1:A:21:PRO:HD2 | 0.48 | 2.39 | 13 | 2 |
| 1:A:18:PHE:CE1 | 1:A:23:THR:HG21 | 0.47 | 2.44 | 14 | 1 |
| 1:A:27:HIS:O | 1:A:31:TYR:HB2 | 0.46 | 2.11 | 19 | 1 |
| 1:A:18:PHE:CE1 | 1:A:23:THR:CG2 | 0.45 | 2.99 | 14 | 2 |
| 1:A:2:SER:O | 1:A:3:LEU:HD23 | 0.45 | 2.11 | 5 | 1 |
| 1:A:18:PHE:CD1 | 1:A:23:THR:CG2 | 0.45 | 2.99 | 1 | 1 |
| 1:A:36:ILE:OXT | 1:A:36:ILE:HG23 | 0.44 | 2.12 | 8 | 1 |
| 1:A:9:PHE:CG | 1:A:9:PHE:O | 0.43 | 2.71 | 16 | 1 |
| 1:A:18:PHE:HB2 | 1:A:24:LEU:CD2 | 0.43 | 2.44 | 13 | 5 |
| 1:A:24:LEU:N | 1:A:24:LEU:HD23 | 0.43 | 2.29 | 10 | 1 |
| 1:A:20:SER:O | 1:A:22:SER:N | 0.43 | 2.52 | 10 | 1 |
| 1:A:16:ILE:HG22 | 1:A:17:ALA:N | 0.43 | 2.29 | 11 | 2 |
| 1:A:32:CYS:O | 1:A:34:HIS:N | 0.43 | 2.52 | 9 | 1 |
| 1:A:10:MET:HG3 | 1:A:17:ALA:HB2 | 0.43 | 1.89 | 6 | 3 |
| 1:A:26:ALA:O | 1:A:31:TYR:CD1 | 0.43 | 2.72 | 9 | 1 |
| 1:A:9:PHE:CD1 | 1:A:9:PHE:O | 0.43 | 2.72 | 16 | 1 |
| 1:A:18:PHE:CD1 | 1:A:23:THR:HG22 | 0.43 | 2.49 | 1 | 1 |
| 1:A:27:HIS:O | 1:A:31:TYR:N | 0.42 | 2.46 | 12 | 1 |
| 1:A:18:PHE:CD1 | 1:A:23:THR:HG21 | 0.42 | 2.50 | 17 | 1 |
| 1:A:18:PHE:CB | 1:A:24:LEU:CD2 | 0.42 | 2.98 | 11 | 2 |
| 1:A:3:LEU:HD12 | 1:A:3:LEU:O | 0.41 | 2.16 | 9 | 1 |
| 1:A:10:MET:CG | 1:A:17:ALA:HB2 | 0.41 | 2.46 | 13 | 1 |
| 1:A:2:SER:OG | 1:A:3:LEU:HD23 | 0.40 | 2.15 | 14 | 1 |
| 1:A:27:HIS:ND1 | 1:A:31:TYR:CB | 0.40 | 2.85 | 19 | 1 |
| 1:A:9:PHE:N | 1:A:9:PHE:CD1 | 0.40 | 2.89 | 4 | 1 |
| 1:A:9:PHE:C | 1:A:9:PHE:CD1 | 0.40 | 2.95 | 16 | 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 | 34/36 (94%) | 23±2 (69±7%) | 8±2 (23±6%) | 3±1 (8±3%) | 2 | 14 |
| All | All | 680/720 (94%) | 468 (69%) | 157 (23%) | 55 (8%) | 2 | 14 |

All 14 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 | 2 | SER | 13 |
| 1 | A | 33 | SER | 11 |
| 1 | A | 4 | LEU | 7 |
| 1 | A | 7 | ALA | 6 |
| 1 | A | 6 | PRO | 4 |
| 1 | A | 3 | LEU | 2 |
| 1 | A | 5 | LYS | 2 |
| 1 | A | 10 | MET | 2 |
| 1 | A | 32 | CYS | 2 |
| 1 | A | 21 | PRO | 2 |
| 1 | A | 22 | SER | 1 |
| 1 | A | 8 | ARG | 1 |
| 1 | A | 14 | CYS | 1 |
| 1 | A | 19 | SER | 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 | 30/30 (100%) | 19±2 (65±7%) | 11±2 (36±7%) | 1 9 |
| All | All | 600/600 (100%) | 387 (64%) | 213 (36%) | 1 9 |

All 22 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 | 3 | LEU | 18 |
| 1 | A | 4 | LEU | 18 |
| 1 | A | 12 | LEU | 16 |
| 1 | A | 28 | GLN | 16 |
| 1 | A | 20 | SER | 14 |
| 1 | A | 31 | TYR | 13 |
| 1 | A | 33 | SER | 11 |
| 1 | A | 5 | LYS | 11 |
| 1 | A | 32 | CYS | 11 |
| 1 | A | 9 | PHE | 10 |

*Continued on next page...*

*Continued from previous page...*

| Mol | Chain | Res | Type | Models (Total) |
|---|---|---|---|---|
| 1 | A | 25 | GLU | 10 |
| 1 | A | 2 | SER | 9 |
| 1 | A | 10 | MET | 8 |
| 1 | A | 8 | ARG | 8 |
| 1 | A | 19 | SER | 8 |
| 1 | A | 23 | THR | 7 |
| 1 | A | 22 | SER | 7 |
| 1 | A | 14 | CYS | 5 |
| 1 | A | 11 | CYS | 5 |
| 1 | A | 35 | ARG | 4 |
| 1 | A | 36 | ILE | 3 |
| 1 | A | 30 | TYR | 1 |

### 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)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

## 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 50% for the well-defined parts and 49% for the entire structure.

## 7.1 Chemical shift list 1

File name: BMRB entry 4644

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 | 247 |
| Number of shifts mapped to atoms | 247 |
| 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 |
|---|---|---|---|
| $^{13}C_{\alpha}$ | 0 | — | — |
| $^{13}C_{\beta}$ | 0 | — | — |
| $^{13}C'$ | 0 | — | — |
| $^{15}N$ | 31 | 0.79 ± 0.64 | None needed (imprecise) |

### 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 50%, i.e. 212 atoms were assigned a chemical shift out of a possible 427. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

| | Total | $^{1}H$ | $^{13}C$ | $^{15}N$ |
|---|---|---|---|---|
| Backbone | 97/169 (57%) | 66/67 (99%) | 0/70 (0%) | 31/32 (97%) |
| Sidechain | 94/209 (45%) | 93/127 (73%) | 0/74 (0%) | 1/8 (12%) |

*Continued on next page...*

*Continued from previous page...*

| | Total | $^{1}$H | $^{13}$C | $^{15}$N |
|---|---|---|---|---|
| Aromatic | 21/49 (43%) | 21/26 (81%) | 0/20 (0%) | 0/3 (0%) |
| Overall | 212/427 (50%) | 180/220 (82%) | 0/164 (0%) | 32/43 (74%) |

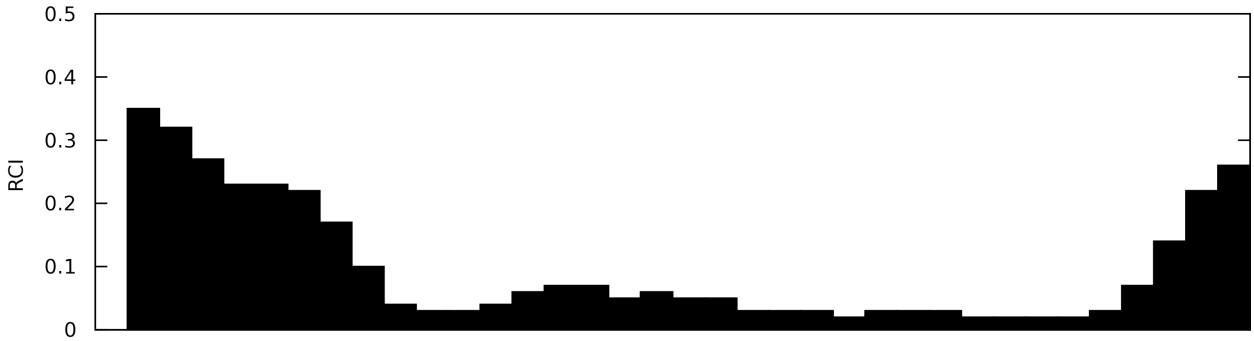
The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 49%, i.e. 212 atoms were assigned a chemical shift out of a possible 432. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

| | Total | $^{1}$H | $^{13}$C | $^{15}$N |
|---|---|---|---|---|
| Backbone | 97/174 (56%) | 66/69 (96%) | 0/72 (0%) | 31/33 (94%) |
| Sidechain | 94/209 (45%) | 93/127 (73%) | 0/74 (0%) | 1/8 (12%) |
| Aromatic | 21/49 (43%) | 21/26 (81%) | 0/20 (0%) | 0/3 (0%) |
| Overall | 212/432 (49%) | 180/222 (81%) | 0/166 (0%) | 32/44 (73%) |

### 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: