



# Full wwPDB NMR Structure Validation Report ⓘ

May 28, 2020 – 11:26 pm BST

PDB ID : 2M3A  
Title : NMR solution structure of a MYB-like DNA binding domain of KNL-2 from C. Elegans  
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Deposited on : 2013-01-15

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

Cyrange : Kirchner and Güntert (2011)  
NmrClust : Kelley et al. (1996)  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
RCI : v\_1n\_11\_5\_13\_A (Berjanski et al., 2005)  
PANAV : Wang et al. (2010)  
ShiftChecker : 2.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

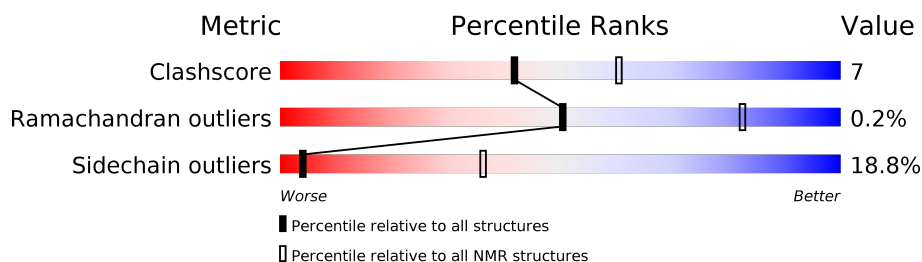
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 53%.

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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  $\geq 3$ , 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  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	67	

## 2 Ensemble composition and analysis

This entry contains 10 models. Model 6 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:10-A:67 (58)	0.25	6

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 1 clusters and 1 single-model cluster was found.

Cluster number	Models
1	1, 2, 3, 4, 5, 6, 8, 9, 10
Single-model clusters	7

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1117 atoms, of which 581 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Protein KNL-2.

Mol	Chain	Residues	Atoms					Trace
1	A	67	Total	C	H	N	O	0
			1117	345	581	98	93	

There are 5 discrepancies between the modelled and reference sequences:

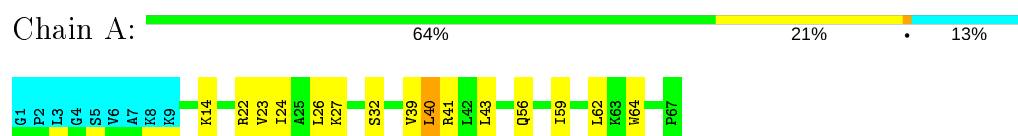
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	EXPRESSION TAG	UNP O44548
A	2	PRO	-	EXPRESSION TAG	UNP O44548
A	3	LEU	-	EXPRESSION TAG	UNP O44548
A	4	GLY	-	EXPRESSION TAG	UNP O44548
A	5	SER	-	EXPRESSION TAG	UNP O44548

## 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: Protein KNL-2

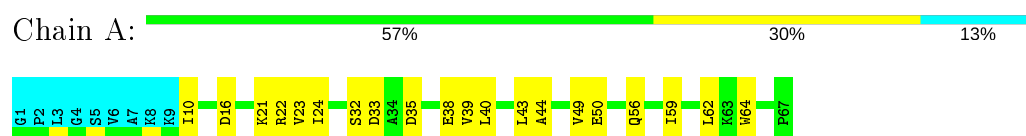


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

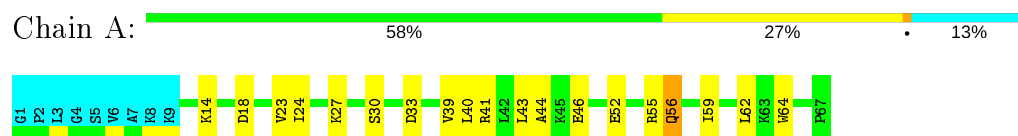
#### 4.2.1 Score per residue for model 1

- Molecule 1: Protein KNL-2



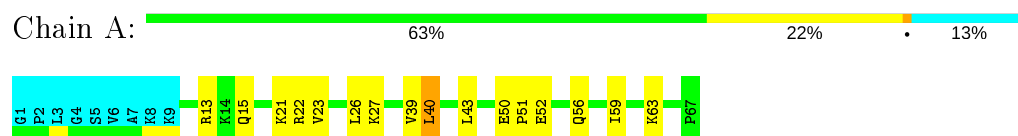
#### 4.2.2 Score per residue for model 2

- Molecule 1: Protein KNL-2



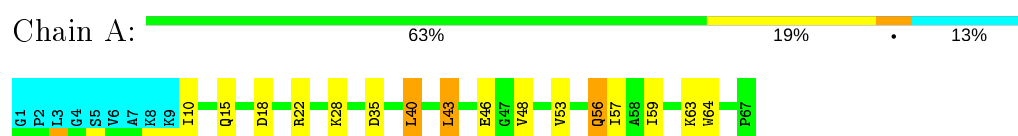
### 4.2.3 Score per residue for model 3

- Molecule 1: Protein KNL-2



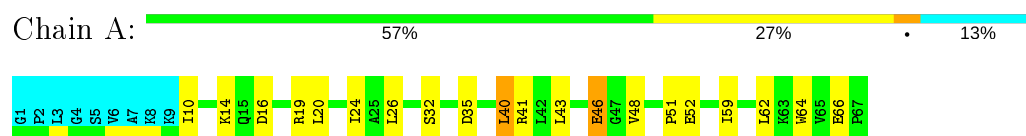
### 4.2.4 Score per residue for model 4

- Molecule 1: Protein KNL-2



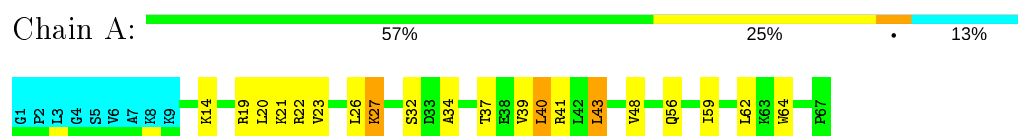
### 4.2.5 Score per residue for model 5

- Molecule 1: Protein KNL-2



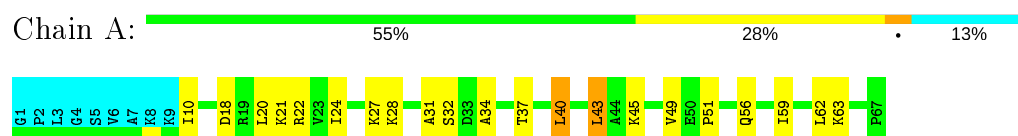
### 4.2.6 Score per residue for model 6 (medoid)

- Molecule 1: Protein KNL-2



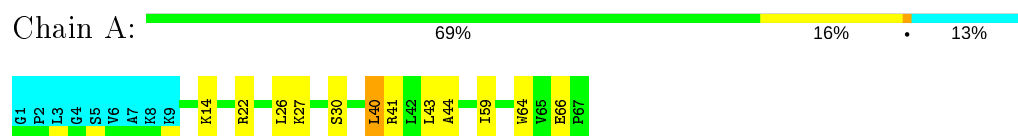
### 4.2.7 Score per residue for model 7

- Molecule 1: Protein KNL-2



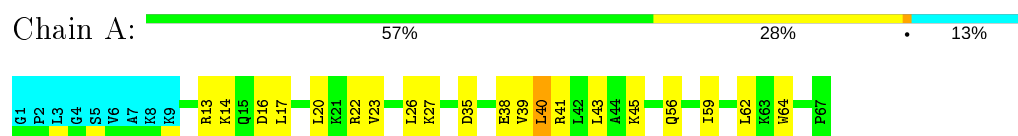
### 4.2.8 Score per residue for model 8

- Molecule 1: Protein KNL-2



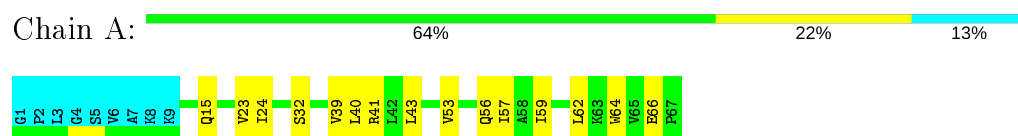
### 4.2.9 Score per residue for model 9

- Molecule 1: Protein KNL-2



### 4.2.10 Score per residue for model 10

- Molecule 1: Protein KNL-2



## 5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 10 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	refinement	

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)	input_cs.cif
Number of chemical shift lists	1
Total number of shifts	568
Number of shifts mapped to atoms	568
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	53%

No validations of the models with respect to experimental NMR restraints is performed at this time.



## 6 Model quality

### 6.1 Standard geometry

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

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	477	512	512	7±2
All	All	4770	5120	5120	72

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 7.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:20:LEU:HD13	1:A:43:LEU:HD12	0.76	1.57	7	2
1:A:24:ILE:HD12	1:A:62:LEU:HD13	0.64	1.67	10	4
1:A:59:ILE:HG23	1:A:64:TRP:O	0.62	1.94	6	8
1:A:23:VAL:HG12	1:A:39:VAL:HG13	0.60	1.71	3	2
1:A:10:ILE:CD1	1:A:49:VAL:HG13	0.59	2.27	1	2
1:A:23:VAL:CG1	1:A:39:VAL:HG13	0.59	2.28	9	6
1:A:56:GLN:HA	1:A:59:ILE:HD12	0.58	1.73	2	7
1:A:21:LYS:HB2	1:A:62:LEU:HD11	0.57	1.75	7	1
1:A:43:LEU:HD22	1:A:48:VAL:CG2	0.57	2.29	6	1
1:A:43:LEU:HD22	1:A:48:VAL:HG21	0.57	1.76	6	2
1:A:46:GLU:HG3	1:A:48:VAL:HG23	0.56	1.77	5	1
1:A:46:GLU:CG	1:A:48:VAL:HG23	0.52	2.35	5	1
1:A:40:LEU:O	1:A:40:LEU:HD12	0.51	2.06	7	3
1:A:24:ILE:HD12	1:A:62:LEU:CD1	0.50	2.36	2	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:40:LEU:HD23	1:A:51:PRO:HB3	0.50	1.84	3	3
1:A:21:LYS:HB2	1:A:62:LEU:HD21	0.49	1.85	1	2
1:A:23:VAL:HB	1:A:39:VAL:HG13	0.48	1.85	2	1
1:A:53:VAL:HG12	1:A:57:ILE:CG1	0.48	2.38	4	2
1:A:20:LEU:CD1	1:A:43:LEU:HD12	0.48	2.33	7	1
1:A:13:ARG:O	1:A:17:LEU:HD13	0.47	2.10	9	1
1:A:20:LEU:HD13	1:A:43:LEU:CD1	0.47	2.37	7	1
1:A:10:ILE:HD12	1:A:49:VAL:HG13	0.47	1.85	7	1
1:A:40:LEU:HD12	1:A:40:LEU:O	0.47	2.09	3	3
1:A:34:ALA:O	1:A:37:THR:HG22	0.46	2.11	7	2
1:A:27:LYS:CB	1:A:39:VAL:HG22	0.44	2.43	6	1
1:A:43:LEU:HD22	1:A:46:GLU:HG2	0.44	1.89	5	1
1:A:43:LEU:HB3	1:A:48:VAL:HG21	0.43	1.90	5	1
1:A:40:LEU:HD21	1:A:51:PRO:HD3	0.43	1.90	5	1
1:A:23:VAL:HG11	1:A:39:VAL:HG13	0.43	1.90	10	1
1:A:23:VAL:CB	1:A:39:VAL:HG13	0.42	2.44	2	1
1:A:40:LEU:O	1:A:44:ALA:HB2	0.42	2.14	8	3
1:A:20:LEU:HG	1:A:62:LEU:HD11	0.42	1.91	9	1
1:A:53:VAL:HG12	1:A:57:ILE:HG12	0.41	1.93	10	1
1:A:21:LYS:CB	1:A:62:LEU:HD11	0.40	2.45	7	1

## 6.3 Torsion angles ⓘ

### 6.3.1 Protein backbone ⓘ

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	57/67 (85%)	51±1 (89±2%)	6±1 (11±1%)	0±0 (0±1%)	50	82
All	All	570/670 (85%)	509 (89%)	60 (11%)	1 (0%)	50	82

All 1 unique Ramachandran outliers are listed below.

Mol	Chain	Res	Type	Models (Total)
1	A	31	ALA	1

### 6.3.2 Protein sidechains ⓘ

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	52/58 (90%)	42±2 (81±4%)	10±2 (19±4%)	4	36
All	All	520/580 (90%)	422 (81%)	98 (19%)	4	36

All 29 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	43	LEU	9
1	A	40	LEU	8
1	A	22	ARG	7
1	A	27	LYS	6
1	A	41	ARG	6
1	A	14	LYS	5
1	A	32	SER	5
1	A	26	LEU	5
1	A	35	ASP	4
1	A	63	LYS	3
1	A	66	GLU	3
1	A	56	GLN	3
1	A	16	ASP	3
1	A	52	GLU	3
1	A	15	GLN	3
1	A	46	GLU	3
1	A	18	ASP	3
1	A	38	GLU	2
1	A	45	LYS	2
1	A	30	SER	2
1	A	19	ARG	2
1	A	33	ASP	2
1	A	28	LYS	2
1	A	50	GLU	2
1	A	20	LEU	1
1	A	55	ARG	1
1	A	13	ARG	1
1	A	10	ILE	1
1	A	21	LYS	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](#)

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 53% for the well-defined parts and 57% for the entire structure.

### 7.1 Chemical shift list 1

File name: input\_cs.cif

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	568
Number of shifts mapped to atoms	568
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	91

#### 7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	47	$0.00 \pm 0.00$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}_\beta$	44	$0.26 \pm 0.12$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
$^{15}\text{N}$	44	$-1.01 \pm 0.44$	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 53%, i.e. 415 atoms were assigned a chemical shift out of a possible 782. 8 out of 14 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	149/284 (52%)	74/113 (65%)	38/116 (33%)	37/55 (67%)
Sidechain	242/462 (52%)	150/269 (56%)	91/167 (54%)	1/26 (4%)

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	<b>Total</b>	<b><sup>1</sup>H</b>	<b><sup>13</sup>C</b>	<b><sup>15</sup>N</b>
Aromatic	24/36 (67%)	12/18 (67%)	10/15 (67%)	2/3 (67%)
Overall	415/782 (53%)	236/400 (59%)	139/298 (47%)	40/84 (48%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 57%, i.e. 500 atoms were assigned a chemical shift out of a possible 880. 10 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	<b>Total</b>	<b><sup>1</sup>H</b>	<b><sup>13</sup>C</b>	<b><sup>15</sup>N</b>
Backbone	181/327 (55%)	90/130 (69%)	47/134 (35%)	44/63 (70%)
Sidechain	295/517 (57%)	183/302 (61%)	111/187 (59%)	1/28 (4%)
Aromatic	24/36 (67%)	12/18 (67%)	10/15 (67%)	2/3 (67%)
Overall	500/880 (57%)	285/450 (63%)	168/336 (50%)	47/94 (50%)

#### 7.1.4 Statistically unusual chemical shifts ⓘ

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	11	THR	CG2	47.70	27.15 – 15.95	23.3
1	A	21	LYS	CG	51.70	30.67 – 19.17	23.3
1	A	45	LYS	CG	51.31	30.67 – 19.17	22.9
1	A	37	THR	CG2	47.22	27.15 – 15.95	22.9
1	A	14	LYS	CG	51.09	30.67 – 19.17	22.8
1	A	8	LYS	CG	50.97	30.67 – 19.17	22.6
1	A	28	LYS	CG	50.94	30.67 – 19.17	22.6
1	A	9	LYS	CG	50.80	30.67 – 19.17	22.5
1	A	27	LYS	CG	50.70	30.67 – 19.17	22.4
1	A	23	VAL	CG1	48.90	28.40 – 14.60	19.9
1	A	39	VAL	CG1	47.34	28.40 – 14.60	18.7
1	A	24	ILE	CG2	43.90	24.63 – 10.43	18.6
1	A	23	VAL	CG2	50.59	29.20 – 13.40	18.5
1	A	10	ILE	CG2	43.49	24.63 – 10.43	18.3
1	A	6	VAL	CG1	46.65	28.40 – 14.60	18.2
1	A	5	SER	CB	38.01	71.24 – 56.34	-17.3
1	A	24	ILE	CD1	42.30	21.91 – 5.01	17.1
1	A	32	SER	CB	38.56	71.24 – 56.34	-16.9
1	A	40	LEU	CD1	51.92	32.77 – 16.57	16.8
1	A	17	LEU	CD1	51.90	32.77 – 16.57	16.8
1	A	29	PRO	CA	37.12	71.13 – 55.53	-16.8
1	A	2	PRO	CA	37.13	71.13 – 55.53	-16.8

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Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	42	LEU	CD1	51.70	32.77 – 16.57	16.7
1	A	30	SER	CB	39.09	71.24 – 56.34	-16.6
1	A	39	VAL	CG2	47.34	29.20 – 13.40	16.5
1	A	6	VAL	CG2	47.34	29.20 – 13.40	16.5
1	A	37	THR	CB	42.05	78.10 – 61.30	-16.5
1	A	40	LEU	CD2	52.03	32.60 – 15.60	16.4
1	A	3	LEU	CD1	50.99	32.77 – 16.57	16.2
1	A	26	LEU	CD1	50.59	32.77 – 16.57	16.0
1	A	26	LEU	CD2	50.38	32.60 – 15.60	15.5
1	A	43	LEU	CD2	49.91	32.60 – 15.60	15.2
1	A	11	THR	CB	44.25	78.10 – 61.30	-15.1
1	A	20	LEU	CD2	49.84	32.60 – 15.60	15.1
1	A	3	LEU	CD2	49.80	32.60 – 15.60	15.1
1	A	42	LEU	CD2	49.30	32.60 – 15.60	14.8
1	A	17	LEU	CD2	49.30	32.60 – 15.60	14.8
1	A	10	ILE	CD1	38.27	21.91 – 5.01	14.7
1	A	31	ALA	CB	45.51	28.03 – 9.93	14.7
1	A	7	ALA	CB	45.40	28.03 – 9.93	14.6
1	A	25	ALA	CB	44.27	28.03 – 9.93	14.0
1	A	34	ALA	CB	44.22	28.03 – 9.93	13.9
1	A	7	ALA	CA	26.38	63.07 – 43.27	-13.5
1	A	44	ALA	CA	26.72	63.07 – 43.27	-13.4
1	A	44	ALA	CB	42.81	28.03 – 9.93	13.2
1	A	45	LYS	CA	28.84	67.97 – 45.97	-12.8
1	A	31	ALA	CA	27.98	63.07 – 43.27	-12.7
1	A	30	SER	CA	32.06	69.25 – 48.25	-12.7
1	A	5	SER	CA	32.15	69.25 – 48.25	-12.7
1	A	27	LYS	CA	29.20	67.97 – 45.97	-12.6
1	A	32	SER	CA	32.52	69.25 – 48.25	-12.5
1	A	25	ALA	CA	28.66	63.07 – 43.27	-12.4
1	A	3	LEU	CA	29.41	66.36 – 44.96	-12.3
1	A	9	LYS	CA	30.07	67.97 – 45.97	-12.2
1	A	8	LYS	CA	30.07	67.97 – 45.97	-12.2
1	A	13	ARG	CA	28.68	68.35 – 45.25	-12.2
1	A	28	LYS	CA	30.48	67.97 – 45.97	-12.0
1	A	46	GLU	CA	32.24	67.86 – 46.86	-12.0
1	A	34	ALA	CA	29.70	63.07 – 43.27	-11.9
1	A	42	LEU	CA	30.69	66.36 – 44.96	-11.7
1	A	35	ASP	CA	31.09	64.88 – 44.48	-11.6
1	A	16	ASP	CA	31.11	64.88 – 44.48	-11.6
1	A	38	GLU	CA	33.09	67.86 – 46.86	-11.6
1	A	26	LEU	CA	31.00	66.36 – 44.96	-11.5

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Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	18	ASP	CA	31.56	64.88 – 44.48	-11.3
1	A	43	LEU	CA	31.64	66.36 – 44.96	-11.2
1	A	33	ASP	CA	31.93	64.88 – 44.48	-11.2
1	A	17	LEU	CA	31.85	66.36 – 44.96	-11.1
1	A	20	LEU	CA	32.11	66.36 – 44.96	-11.0
1	A	15	GLN	CA	33.15	67.31 – 45.91	-11.0
1	A	40	LEU	CA	32.75	66.36 – 44.96	-10.7
1	A	12	TRP	CA	30.70	70.54 – 44.84	-10.5
1	A	10	ILE	CA	33.85	75.08 – 48.18	-10.3
1	A	11	THR	CA	35.16	75.37 – 49.07	-10.3
1	A	21	LYS	CA	34.34	67.97 – 45.97	-10.3
1	A	22	ARG	CA	33.39	68.35 – 45.25	-10.1
1	A	14	LYS	CA	34.88	67.97 – 45.97	-10.0
1	A	41	ARG	CA	33.61	68.35 – 45.25	-10.0
1	A	19	ARG	CA	33.66	68.35 – 45.25	-10.0
1	A	36	TRP	NE1	110.40	139.19 – 119.59	-9.7
1	A	12	TRP	NE1	111.38	139.19 – 119.59	-9.2
1	A	6	VAL	CA	36.15	76.93 – 48.03	-9.1
1	A	36	TRP	CA	35.26	70.54 – 44.84	-8.7
1	A	36	TRP	HZ2	4.60	8.89 – 5.69	-8.4
1	A	45	LYS	HE2	1.40	3.87 – 1.97	-8.0
1	A	24	ILE	CA	40.23	75.08 – 48.18	-8.0
1	A	37	THR	CA	41.60	75.37 – 49.07	-7.8
1	A	39	VAL	CA	40.33	76.93 – 48.03	-7.7
1	A	23	VAL	CA	40.89	76.93 – 48.03	-7.5
1	A	45	LYS	HE3	1.75	3.86 – 1.96	-6.1
1	A	29	PRO	HG3	0.25	3.56 – 0.26	-5.0

### 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:



