



# Full wwPDB NMR Structure Validation Report ⓘ

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PDB ID : 1K8J  
Title : NMR STRUCTURE OF THE CK14 DNA DUPLEX: A PORTION OF THE KNOWN NF-kB SEQUENCE CK1  
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We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/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	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	trunk28760
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*


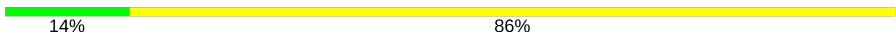
The overall completeness of chemical shifts assignment is 33%.

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	125131	11601

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	14	 14% 86%
2	B	14	 14% 86%

## 2 Ensemble composition and analysis ⓘ

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

### 3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 886 atoms, of which 318 are hydrogens and 0 are deuteriums.

- Molecule 1 is a DNA chain called FIRST STRAND OF CK14 DNA DUPLEX.

Mol	Chain	Residues	Atoms						Trace
1	A	14	Total	C	H	N	O	P	0
			440	135	158	54	80	13	

- Molecule 2 is a DNA chain called SECOND STRAND OF CK14 DNA DUPLEX.

Mol	Chain	Residues	Atoms						Trace
2	B	14	Total	C	H	N	O	P	0
			446	137	160	52	84	13	

## 4 Residue-property plots

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 STRAND OF CK14 DNA DUPLEX

Chain A: 

C1	C2	A3	G4	G5	A6	G7	A8	T9	T10	C11	C12	A13	C14
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- Molecule 2: SECOND STRAND OF CK14 DNA DUPLEX

Chain B: 

G15	T16	G17	G18	A19	A20	T21	C22	C23	C24	C25	T26	G27	G28
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## 5 Refinement protocol and experimental data overview

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

Of the 10 calculated structures, 1 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
AMBER	refinement	5

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 5718, BMRB entry 5717, BMRB entry 5716
Number of chemical shift lists	5
Total number of shifts	598
Number of shifts mapped to atoms	526
Number of unparsed shifts	0
Number of shifts with mapping errors	72
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	33%

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

## 6 Model quality

### 6.1 Standard geometry

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	1.37	0/316 (0.0%)	2.05	23/485 (4.7%)
2	B	1.41	0/320 (0.0%)	2.07	19/493 (3.9%)
All	All	1.39	0/636 (0.0%)	2.06	42/978 (4.3%)

There are no bond-length outliers.

All angle outliers are listed below. They are sorted according to the Z-score.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	17	DG	O4'-C1'-N9	7.32	113.12	108.00
2	B	21	DT	C6-C5-C7	-6.20	119.18	122.90
2	B	27	DG	O4'-C1'-N9	6.17	112.32	108.00
2	B	20	DA	C5-C6-N1	6.01	120.71	117.70
1	A	13	DA	C5-C6-N1	5.97	120.69	117.70
1	A	3	DA	C5-C6-N1	5.96	120.68	117.70
1	A	6	DA	C5-C6-N1	5.93	120.67	117.70
1	A	11	DC	N3-C2-O2	-5.89	117.78	121.90
2	B	22	DC	O4'-C1'-N1	5.81	112.07	108.00
1	A	13	DA	N1-C6-N6	-5.79	115.13	118.60
1	A	8	DA	C5-C6-N1	5.74	120.57	117.70
1	A	9	DT	C6-C5-C7	-5.72	119.47	122.90
2	B	19	DA	C5-C6-N1	5.71	120.55	117.70
1	A	12	DC	N3-C2-O2	-5.62	117.97	121.90
1	A	10	DT	C6-C5-C7	-5.61	119.53	122.90
1	A	6	DA	N1-C6-N6	-5.54	115.28	118.60
1	A	1	DC	N3-C2-O2	-5.53	118.03	121.90
2	B	24	DC	N3-C2-O2	-5.52	118.04	121.90
1	A	11	DC	N1-C2-O2	5.51	122.21	118.90
1	A	3	DA	N1-C6-N6	-5.43	115.34	118.60
2	B	25	DC	N3-C2-O2	-5.41	118.11	121.90
1	A	2	DC	N3-C2-O2	-5.41	118.11	121.90
2	B	24	DC	N1-C2-O2	5.37	122.12	118.90
2	B	27	DG	N1-C6-O6	-5.35	116.69	119.90
1	A	3	DA	C4-C5-C6	-5.32	114.34	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	DA	N1-C6-N6	-5.29	115.42	118.60
2	B	22	DC	N3-C2-O2	-5.29	118.20	121.90
2	B	26	DT	C6-C5-C7	-5.27	119.74	122.90
2	B	19	DA	N1-C6-N6	-5.26	115.44	118.60
2	B	23	DT	C6-C5-C7	-5.22	119.77	122.90
2	B	28	DG	N1-C6-O6	-5.20	116.78	119.90
1	A	12	DC	N1-C2-O2	5.18	122.01	118.90
2	B	16	DT	C6-C5-C7	-5.14	119.82	122.90
1	A	6	DA	C4-C5-C6	-5.09	114.46	117.00
1	A	14	DC	N3-C2-O2	-5.09	118.34	121.90
1	A	8	DA	C4-C5-C6	-5.08	114.46	117.00
1	A	4	DG	N1-C6-O6	-5.06	116.86	119.90
1	A	1	DC	N1-C2-O2	5.06	121.94	118.90
2	B	19	DA	C4-C5-C6	-5.05	114.47	117.00
2	B	21	DT	N3-C2-O2	-5.03	119.28	122.30
2	B	25	DC	N1-C2-O2	5.01	121.91	118.90
1	A	10	DT	N3-C2-O2	-5.00	119.30	122.30

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
All	All	568	318	318	-

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 -.

There are no clashes.

## 6.3 Torsion angles [i](#)

### 6.3.1 Protein backbone [i](#)

There are no protein molecules in this entry.



### 6.3.2 Protein sidechains [i](#)

There are no protein molecules in this entry.

### 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

The completeness of assignment taking into account all chemical shift lists is 33% for the well-defined parts and 33% for the entire structure.

### 7.1 Chemical shift list 1

File name: BMRB entry 5716

Chemical shift list name: *assigned\_chem\_shift\_list\_1*

#### 7.1.1 Bookkeeping

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	99
Number of shifts mapped to atoms	99
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

No chemical shift referencing corrections were calculated (not enough data).

#### 7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 17%, i.e. 93 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 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	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	93/556 (17%)	93/332 (28%)	0/188 (0%)	0/36 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 17%, i.e. 93 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	93/556 (17%)	93/332 (28%)	0/188 (0%)	0/36 (0%)

### 7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

### 7.1.5 Random Coil Index (RCI) plots [i](#)

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned\_chem\_shift\_list\_1). RCI is only applicable to proteins.

## 7.2 Chemical shift list 2

File name: BMRB entry 5717

Chemical shift list name: *assigned\_chem\_shift\_list\_1*

### 7.2.1 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	200
Number of shifts mapped to atoms	128
Number of unparsed shifts	0
Number of shifts with mapping errors	72
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- Residue not found in structure. All 72 occurrences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
A	10	DC	H6	7.71	0.01	1
A	10	DC	H5	5.68	0.01	1
A	1	DG	H2'	2.6	0.01	1

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Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
A	5	DA	H8	8.05	0.01	1
A	13	DG	H8	8.14	0.01	1
A	12	DT	H73	1.7	0.01	1
A	7	DT	H72	1.3	0.01	1
A	5	DA	H2''	2.89	0.01	1
A	8	DC	H2'	2.17	0.01	1
A	1	DG	H4'	4.28	0.01	1
A	2	DT	H71	1.41	0.01	1
A	10	DC	H1'	6.01	0.01	1
A	12	DT	H6	7.41	0.01	1
A	7	DT	H1'	5.69	0.01	1
A	8	DC	H6	7.67	0.01	1
A	8	DC	H5	5.52	0.01	1
A	3	DG	H1'	5.7	0.01	1
A	12	DT	H2''	2.27	0.01	1
A	3	DG	H2''	2.76	0.01	1
A	8	DC	H2''	2.56	0.01	1
A	14	DG	H2'	2.57	0.01	1
A	13	DG	H2'	2.73	0.01	1
A	7	DT	H4'	4.28	0.01	1
A	5	DA	H2'	2.66	0.01	1
A	2	DT	H2'	2.15	0.01	1
A	2	DT	H6	7.35	0.01	1
A	5	DA	H1'	5.94	0.01	1
A	12	DT	H3'	5.03	0.01	1
A	5	DA	H2	7.29	0.01	1
A	10	DC	H2''	2.51	0.01	1
A	2	DT	H2''	2.43	0.01	1
A	7	DT	H3'	5.05	0.01	1
A	2	DT	H3'	5.09	0.01	1
A	5	DA	H3'	5.04	0.01	1
A	3	DG	H2'	2.65	0.01	1
A	12	DT	H71	1.7	0.01	1
A	14	DG	H3'	4.68	0.01	1
A	7	DT	H6	7.12	0.01	1
A	14	DG	H4'	4.2	0.01	1
A	8	DC	H1'	5.95	0.01	1
A	2	DT	H72	1.41	0.01	1
A	14	DG	H8	7.83	0.01	1
A	3	DG	H3'	4.99	0.01	1
A	7	DT	H73	1.3	0.01	1

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Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
A	1	DG	H1'	6.0	0.01	1
A	7	DT	H2'	2.17	0.01	1
A	8	DC	H3'	4.68	0.01	1
A	14	DG	H1'	6.13	0.01	1
A	1	DG	H2''	2.77	0.01	1
A	1	DG	H3'	5.13	0.01	1
A	10	DC	H3'	4.83	0.01	1
A	12	DT	H4'	4.15	0.01	1
A	8	DC	H4'	4.26	0.01	1
A	3	DG	H4'	4.4	0.01	1
A	7	DT	H2''	2.43	0.01	1
A	12	DT	H72	1.7	0.01	1
A	12	DT	H2'	2.08	0.01	1
A	13	DG	H1'	5.79	0.01	1
A	13	DG	H2''	2.78	0.01	1
A	2	DT	H73	1.41	0.01	1
A	10	DC	H4'	4.23	0.01	1
A	2	DT	H1'	5.52	0.01	1
A	2	DT	H4'	4.23	0.01	1
A	7	DT	H71	1.3	0.01	1
A	13	DG	H3'	5.0	0.01	1
A	14	DG	H2''	2.35	0.01	1
A	5	DA	H4'	4.43	0.01	1
A	13	DG	H4'	4.42	0.01	1
A	3	DG	H8	8.02	0.01	1
A	1	DG	H8	7.95	0.01	1
A	12	DT	H1'	5.31	0.01	1
A	10	DC	H2'	2.23	0.01	1

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

No chemical shift referencing corrections were calculated (not enough data).

## 7.2.3 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 20%, i.e. 113 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	113/556 (20%)	113/332 (34%)	0/188 (0%)	0/36 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 20%, i.e. 113 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	113/556 (20%)	113/332 (34%)	0/188 (0%)	0/36 (0%)

#### 7.2.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

#### 7.2.5 Random Coil Index (RCI) plots [i](#)

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned\_chem\_shift\_list\_1). RCI is only applicable to proteins.

### 7.3 Chemical shift list 3

File name: BMRB entry 5718

Chemical shift list name: *assigned\_chem\_shift\_list\_1*

#### 7.3.1 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	97
Number of shifts mapped to atoms	97
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

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

No chemical shift referencing corrections were calculated (not enough data).

### 7.3.3 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 16%, i.e. 91 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	91/556 (16%)	91/332 (27%)	0/188 (0%)	0/36 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 16%, i.e. 91 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	91/556 (16%)	91/332 (27%)	0/188 (0%)	0/36 (0%)

### 7.3.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

### 7.3.5 Random Coil Index (RCI) plots [i](#)

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned\_chem\_shift\_list\_1). RCI is only applicable to proteins.

## 7.4 Chemical shift list 4

File name: BMRB entry 5716

Chemical shift list name: *assigned\_chem\_shift\_list\_2*

### 7.4.1 Bookkeeping

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	101
Number of shifts mapped to atoms	101
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

### 7.4.2 Chemical shift referencing

No chemical shift referencing corrections were calculated (not enough data).

### 7.4.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 16%, i.e. 89 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 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	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	89/556 (16%)	89/332 (27%)	0/188 (0%)	0/36 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 16%, i.e. 89 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 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	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	89/556 (16%)	89/332 (27%)	0/188 (0%)	0/36 (0%)

### 7.4.4 Statistically unusual chemical shifts

There are no statistically unusual chemical shifts.



### 7.4.5 Random Coil Index (RCI) plots [i](#)

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned\_chem\_shift\_list\_2). RCI is only applicable to proteins.

## 7.5 Chemical shift list 5

File name: BMRB entry 5718

Chemical shift list name: *assigned\_chem\_shift\_list\_2*

### 7.5.1 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	101
Number of shifts mapped to atoms	101
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

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

No chemical shift referencing corrections were calculated (not enough data).

### 7.5.3 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 16%, i.e. 89 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 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	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	89/556 (16%)	89/332 (27%)	0/188 (0%)	0/36 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 16%, i.e. 89 atoms were assigned a chemical shift out of a possible 556. 0 out of 0 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	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Sidechain	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Aromatic	0/0 (—%)	0/0 (—%)	0/0 (—%)	0/0 (—%)
Overall	89/556 (16%)	89/332 (27%)	0/188 (0%)	0/36 (0%)

#### 7.5.4 Statistically unusual chemical shifts ⓘ

There are no statistically unusual chemical shifts.

#### 7.5.5 Random Coil Index (RCI) plots ⓘ

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned\_chem\_shift\_list\_2). RCI is only applicable to proteins.