



Full wwPDB NMR Structure Validation Report ⓘ

Feb 12, 2017 – 06:20 pm GMT

PDB ID : 1I5T
Title : SOLUTION STRUCTURE OF CYANOFERRICYTOCHROME C
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Deposited on : 2001-02-28

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A user guide is available at

<http://wwpdb.org/validation/2016/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
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

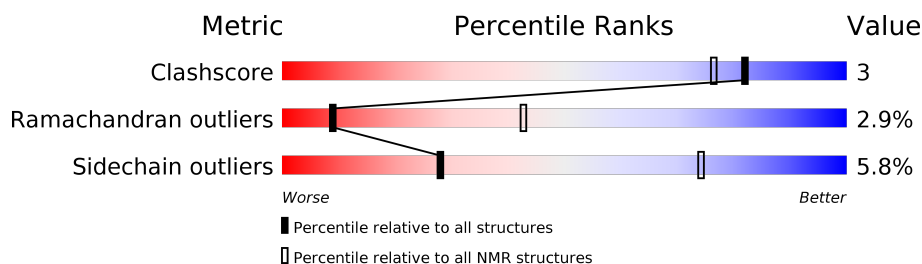
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 44%.

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
Ramachandran outliers	121729	10391
Sidechain outliers	121581	10367

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	104	 84% 14% .

2 Ensemble composition and analysis ⓘ

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

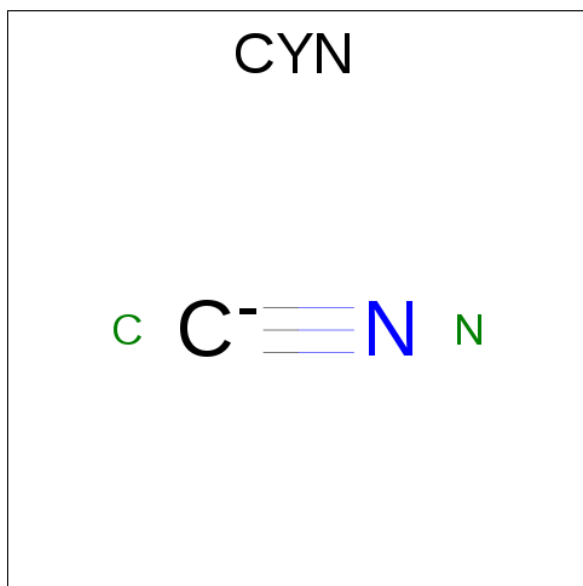
3 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1747 atoms, of which 879 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called CYTOCHROME C.

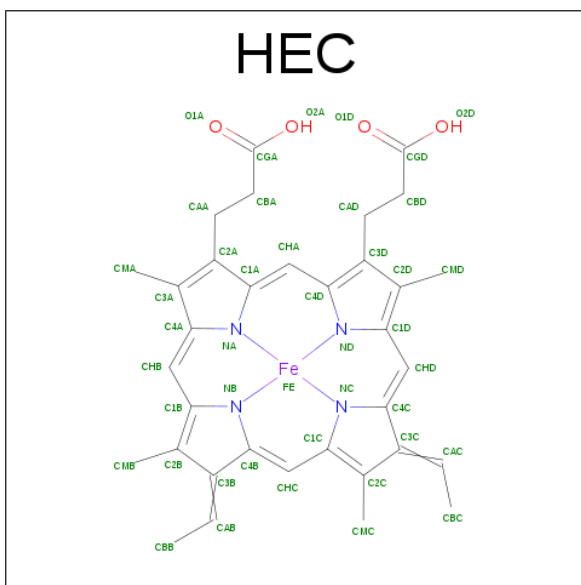
Mol	Chain	Residues	Atoms						Trace
1	A	104	Total	C	H	N	O	S	0
			1670	524	847	144	151	4	

- Molecule 2 is CYANIDE ION (three-letter code: CYN) (formula: CN).



Mol	Chain	Residues	Atoms		
2	A	1	Total	C	N
			2	1	1

- Molecule 3 is HEME C (three-letter code: HEC) (formula: C₃₄H₃₄FeN₄O₄).




Mol	Chain	Residues	Atoms					
3	A	1	Total	C	Fe	H	N	O
			75	34	1	32	4	4

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: CYTOCHROME C

Chain A:  84% 14% .



5 Refinement protocol and experimental data overview

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

Of the 35 calculated structures, 1 were deposited, based on the following criterion: *all calculated structures submitted*.

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

Software name	Classification	Version
DYANA	structure solution	1.5
AMBER	refinement	5.0

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 5026
Number of chemical shift lists	2
Total number of shifts	722
Number of shifts mapped to atoms	690
Number of unparsed shifts	0
Number of shifts with mapping errors	32
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	44%

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

6 Model quality

6.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CYN, HEC

There are no covalent bond-length or bond-angle outliers.

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	6
All	All	0	6

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

All planar outliers are listed below.

Mol	Chain	Res	Type	Group
1	A	48	TYR	Sidechain
1	A	67	TYR	Sidechain
1	A	91	ARG	Sidechain
1	A	18	HIS	Sidechain
1	A	78	THR	Peptide
1	A	74	TYR	Sidechain

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	823	847	847	4
3	A	43	32	30	1
All	All	868	879	877	5

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 3.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:9:ILE:HG23	1:A:13:LYS:HE3	0.56	1.76
1:A:9:ILE:HG23	1:A:13:LYS:CE	0.54	2.33
1:A:97:TYR:CE1	1:A:101:ALA:HB2	0.53	2.39
3:A:110:HEC:HMC1	3:A:110:HEC:HBC3	0.49	1.84
1:A:20:VAL:HG21	1:A:101:ALA:HB1	0.42	1.91

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	102/104 (98%)	84 (82%)	15 (15%)	3 (3%)	9	43
All	All	102/104 (98%)	84 (82%)	15 (15%)	3 (3%)	9	43

All 3 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	80	MET
1	A	33	HIS
1	A	81	ILE

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	86/86 (100%)	81 (94%)	5 (6%)	28	74
All	All	86/86 (100%)	81 (94%)	5 (6%)	28	74

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	26	HIS
1	A	74	TYR
1	A	13	LYS
1	A	3	VAL
1	A	89	THR

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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
2	CYN	A	105	3	0,1,1	0.00	-
3	HEC	A	110	1,2	28,50,50	0.98	0 (0%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard

deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
2	CYN	A	105	3	0,0,0	0.00	-
3	HEC	A	110	1,2	16,82,82	1.22	0 (0%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CYN	A	105	3	-	0,0,0,0	0,0,0,0
3	HEC	A	110	1,2	-	0,6,54,54	0,0,8,8

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 44% for the well-defined parts and 44% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 5026

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	690
Number of shifts mapped to atoms	690
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	37

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 44%, i.e. 580 atoms were assigned a chemical shift out of a possible 1308. 0 out of 9 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	198/512 (39%)	198/204 (97%)	0/208 (0%)	0/100 (0%)
Sidechain	335/695 (48%)	335/411 (82%)	0/251 (0%)	0/33 (0%)
Aromatic	47/101 (47%)	47/54 (87%)	0/43 (0%)	0/4 (0%)
Overall	580/1308 (44%)	580/669 (87%)	0/502 (0%)	0/137 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 44%, i.e. 580 atoms were assigned a chemical shift out of a possible 1308. 0 out of 9 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	198/512 (39%)	198/204 (97%)	0/208 (0%)	0/100 (0%)
Sidechain	335/695 (48%)	335/411 (82%)	0/251 (0%)	0/33 (0%)
Aromatic	47/101 (47%)	47/54 (87%)	0/43 (0%)	0/4 (0%)
Overall	580/1308 (44%)	580/669 (87%)	0/502 (0%)	0/137 (0%)

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	18	HIS	HE1	-11.95	10.53 – 5.43	-39.1
1	A	18	HIS	HD2	22.18	9.28 – 4.78	33.7
1	A	18	HIS	HB2	11.77	4.91 – 1.31	24.1
1	A	18	HIS	HB3	8.71	5.00 – 1.10	14.5
1	A	18	HIS	HA	9.65	6.81 – 2.41	11.5
1	A	67	TYR	HH	22.78	16.26 – 2.26	9.7
1	A	71	PRO	HB3	5.34	3.81 – 0.21	9.2
1	A	71	PRO	HB2	5.25	3.82 – 0.32	9.1
1	A	67	TYR	HE1	8.65	7.86 – 5.56	8.4
1	A	67	TYR	HE2	8.65	7.86 – 5.56	8.4
1	A	72	LYS	HD3	3.46	2.75 – 0.45	8.1
1	A	72	LYS	HD2	3.46	2.76 – 0.46	8.0
1	A	75	ILE	HD12	2.95	2.13 – -0.77	7.8
1	A	75	ILE	HD11	2.95	2.13 – -0.77	7.8
1	A	75	ILE	HD13	2.95	2.13 – -0.77	7.8
1	A	32	LEU	HG	4.09	3.16 – -0.14	7.8
1	A	41	GLY	HA3	1.25	5.80 – 2.00	-7.0
1	A	68	LEU	HD23	-1.14	2.14 – -0.66	-6.7
1	A	68	LEU	HD21	-1.14	2.14 – -0.66	-6.7
1	A	68	LEU	HD22	-1.14	2.14 – -0.66	-6.7
1	A	91	ARG	HD2	1.73	4.27 – 1.97	-6.0
1	A	64	LEU	HD23	-0.94	2.14 – -0.66	-6.0
1	A	64	LEU	HD22	-0.94	2.14 – -0.66	-6.0
1	A	64	LEU	HD21	-0.94	2.14 – -0.66	-6.0
1	A	28	THR	HG21	-0.23	2.29 – -0.01	-6.0
1	A	28	THR	HG23	-0.23	2.29 – -0.01	-6.0
1	A	28	THR	HG22	-0.23	2.29 – -0.01	-6.0
1	A	78	THR	HB	6.09	5.82 – 2.52	5.8
1	A	31	ASN	H	11.91	11.45 – 5.25	5.7
1	A	75	ILE	HB	3.41	3.24 – 0.34	5.6
1	A	91	ARG	HD3	1.73	4.36 – 1.86	-5.5

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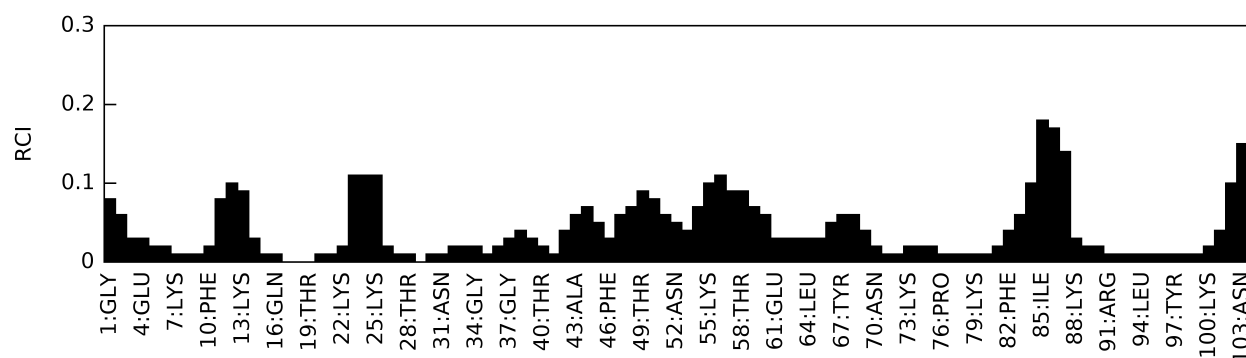
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Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	35	LEU	HD23	-0.72	2.14 – -0.66	-5.2
1	A	35	LEU	HD21	-0.72	2.14 – -0.66	-5.2
1	A	35	LEU	HD22	-0.72	2.14 – -0.66	-5.2
1	A	19	THR	HG22	2.33	2.29 – -0.01	5.2
1	A	19	THR	HG23	2.33	2.29 – -0.01	5.2
1	A	19	THR	HG21	2.33	2.29 – -0.01	5.2

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:



7.2 Chemical shift list 2

File name: BMRB entry 5026

Chemical shift list name: *assigned_chem_shift_list_2*

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	32
Number of shifts mapped to atoms	0
Number of unparsed shifts	0
Number of shifts with mapping errors	32

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.

- Chain not found in structure. All 32 occurrences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
UNMAPPED	1	HEC	HHC	8.99	-1.0	1
UNMAPPED	1	HEC	HMB1	16.49	-1.0	1
UNMAPPED	1	HEC	HMD2	22.96	-1.0	1
UNMAPPED	1	HEC	HBB1	0.42	-1.0	1
UNMAPPED	1	HEC	HMD1	22.96	-1.0	1
UNMAPPED	1	HEC	HMA1	21.48	-1.0	1
UNMAPPED	1	HEC	HBD1	2.09	-1.0	1
UNMAPPED	1	HEC	HMA2	21.48	-1.0	1
UNMAPPED	1	HEC	HBD2	2.81	-1.0	1
UNMAPPED	1	HEC	HMC1	11.4	-1.0	1
UNMAPPED	1	HEC	HBC3	0.23	-1.0	1
UNMAPPED	1	HEC	HAD2	5.01	-1.0	1
UNMAPPED	1	HEC	HAD1	3.12	-1.0	1
UNMAPPED	1	HEC	HHB	0.59	-1.0	1
UNMAPPED	1	HEC	HMB3	16.49	-1.0	1
UNMAPPED	1	HEC	HBB2	0.42	-1.0	1
UNMAPPED	1	HEC	HMA3	21.48	-1.0	1
UNMAPPED	1	HEC	HMC2	11.4	-1.0	1
UNMAPPED	1	HEC	HBC2	0.23	-1.0	1
UNMAPPED	1	HEC	HAB	0.54	-1.0	1
UNMAPPED	1	HEC	HHA	5.71	-1.0	1
UNMAPPED	1	HEC	HAA1	2.14	-1.0	1
UNMAPPED	1	HEC	HHH	2.81	-1.0	1
UNMAPPED	1	HEC	HAA2	5.49	-1.0	1
UNMAPPED	1	HEC	HMB2	16.49	-1.0	1
UNMAPPED	1	HEC	HBB3	0.42	-1.0	1
UNMAPPED	1	HEC	HMD3	22.96	-1.0	1
UNMAPPED	1	HEC	HBA2	0.22	-1.0	1
UNMAPPED	1	HEC	HBA1	-0.1	-1.0	1
UNMAPPED	1	HEC	HMC3	11.4	-1.0	1
UNMAPPED	1	HEC	HBC1	0.23	-1.0	1
UNMAPPED	1	HEC	HAC	-0.38	-1.0	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 0%, i.e. 0 atoms were assigned a chemical shift out of a possible 1308. 0 out of 9 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/512 (0%)	0/204 (0%)	0/208 (0%)	0/100 (0%)
Sidechain	0/695 (0%)	0/411 (0%)	0/251 (0%)	0/33 (0%)
Aromatic	0/101 (0%)	0/54 (0%)	0/43 (0%)	0/4 (0%)
Overall	0/1308 (0%)	0/669 (0%)	0/502 (0%)	0/137 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 0%, i.e. 0 atoms were assigned a chemical shift out of a possible 1308. 0 out of 9 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/512 (0%)	0/204 (0%)	0/208 (0%)	0/100 (0%)
Sidechain	0/695 (0%)	0/411 (0%)	0/251 (0%)	0/33 (0%)
Aromatic	0/101 (0%)	0/54 (0%)	0/43 (0%)	0/4 (0%)
Overall	0/1308 (0%)	0/669 (0%)	0/502 (0%)	0/137 (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_2). RCI is only applicable to proteins.