



wwPDB NMR Structure Validation Summary Report ⓘ

Apr 26, 2016 – 02:12 PM BST

PDB ID : 1EHJ
Title : A PROTON-NMR INVESTIGATION OF THE FULLY REDUCED CY-
TOCHROME C7 FROM DESULFUROMONAS ACETOXIDANS
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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 ⓘ 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 : 1.7.1 (RC1), CSD as537be (2016)
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

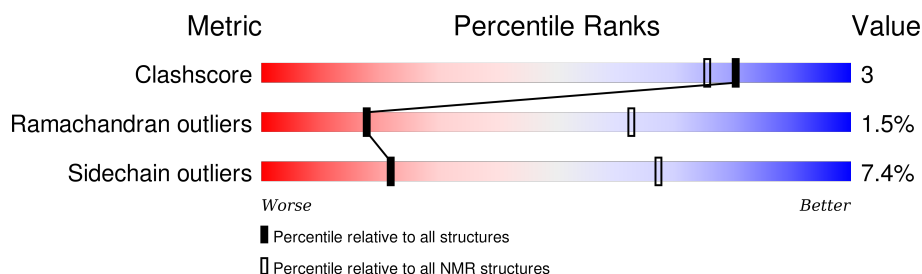
1 Overall quality at a glance

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	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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	68	

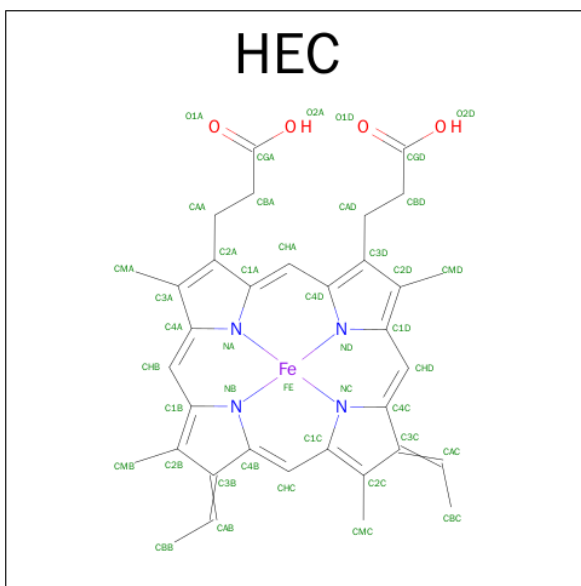
2 Ensemble composition and analysis ⓘ

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

- Molecule 1 is a protein called CYTOCHROME C7.

Mol	Chain	Residues	Atoms						Trace
1	A	68	Total	C	H	N	O	S	0
			997	306	492	96	97	6	

- Molecule 2 is HEME C (three-letter code: HEC) (formula: $\text{C}_{34}\text{H}_{34}\text{FeN}_4\text{O}_4$).



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

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: CYTOCHROME C7

Chain A:  90% 10%



5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing combined with torsion angle dynamics restrained energy minimization*.

Of the ? calculated structures, 1 were deposited, based on the following criterion: ?.

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

Software name	Classification	Version
AMBER	refinement	4.1
DYANA	structure solution	1.5

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 4743
Number of chemical shift lists	4
Total number of shifts	449
Number of shifts mapped to atoms	400
Number of unparsed shifts	0
Number of shifts with mapping errors	49
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 [i](#)

6.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1	0
All	All	1	0

There are no bond-length outliers.

There are no bond-angle outliers.

All chiral outliers are listed below.

Mol	Chain	Res	Type	Atoms
1	A	1	ALA	CA

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	505	492	492	2
2	A	129	96	90	4
All	All	634	588	582	4

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)
2:A:1030:HEC:HBB3	2:A:1030:HEC:HMB1	0.74	1.60
1:A:13:VAL:HG11	2:A:1066:HEC:HMD3	0.59	1.73
2:A:1030:HEC:CBB	2:A:1030:HEC:HMB1	0.45	2.37
1:A:36:LYS:NZ	2:A:1030:HEC:O1D	0.41	2.53

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	66/68 (97%)	39 (59%)	26 (39%)	1 (2%)	18	63
All	All	66/68 (97%)	39 (59%)	26 (39%)	1 (2%)	18	63

All 1 Ramachandran outliers are listed below.

Mol	Chain	Res	Type
1	A	26	CYS

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	54/54 (100%)	50 (93%)	4 (7%)	22	67
All	All	54/54 (100%)	50 (93%)	4 (7%)	22	67

All 4 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	52	CYS
1	A	41	LYS
1	A	51	THR

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Mol	Chain	Res	Type
1	A	68	LYS

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](#)

3 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	HEC	A	1030	1	24,50,50	1.01	0 (0%)
2	HEC	A	1053	1	24,50,50	0.97	0 (0%)
2	HEC	A	1066	1	24,50,50	1.01	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	HEC	A	1030	1	19,82,82	2.05	1 (5%)
2	HEC	A	1053	1	19,82,82	2.23	2 (10%)
2	HEC	A	1066	1	19,82,82	2.51	2 (10%)

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	HEC	A	1030	1	-	0,6,54,54	0,0,8,8
2	HEC	A	1053	1	-	0,6,54,54	0,0,8,8
2	HEC	A	1066	1	-	0,6,54,54	0,0,8,8

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	A	1053	HEC	CBB-CAB-C3B	6.64	112.83	127.34
2	A	1066	HEC	CBB-CAB-C3B	6.17	113.84	127.34
2	A	1066	HEC	CBC-CAC-C3C	6.13	113.95	127.34
2	A	1030	HEC	CBC-CAC-C3C	5.55	115.21	127.34
2	A	1053	HEC	CBC-CAC-C3C	5.44	115.45	127.34

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 4743

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	400
Number of shifts mapped to atoms	400
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	34

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. 342 atoms were assigned a chemical shift out of a possible 776. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	128/336 (38%)	128/134 (96%)	0/136 (0%)	0/66 (0%)
Sidechain	199/381 (52%)	199/225 (88%)	0/140 (0%)	0/16 (0%)
Aromatic	15/59 (25%)	15/33 (45%)	0/20 (0%)	0/6 (0%)
Overall	342/776 (44%)	342/392 (87%)	0/296 (0%)	0/88 (0%)

7.1.4 Statistically unusual chemical shifts ⓘ

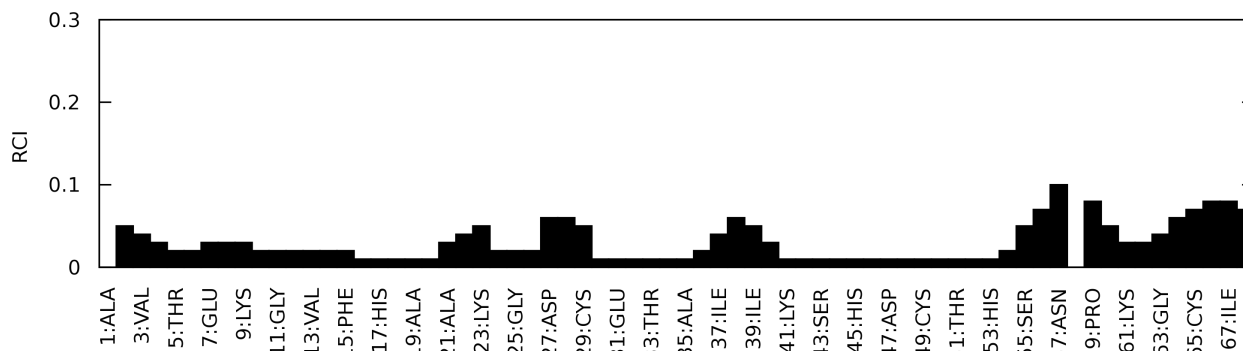
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, <i>ppm</i>	Expected range, <i>ppm</i>	Z-score
1	A	53	HIS	HE1	0.72	10.53 – 5.43	-14.2
1	A	34	PRO	HG2	-2.35	3.48 – 0.38	-13.8
1	A	66	HIS	HE1	1.01	10.53 – 5.43	-13.7
1	A	10	LYS	HE3	0.59	3.86 – 1.96	-12.2
1	A	41	LYS	HD2	-1.06	2.76 – 0.46	-11.6
1	A	10	LYS	HE2	0.72	3.87 – 1.97	-11.6
1	A	10	LYS	HG3	-1.83	2.76 – -0.04	-11.4
1	A	15	PHE	HE2	10.20	8.69 – 5.49	9.7
1	A	15	PHE	HE1	10.20	8.69 – 5.49	9.7
1	A	33	THR	HB	1.12	5.82 – 2.52	-9.2
1	A	41	LYS	HG3	-1.16	2.76 – -0.04	-9.0
1	A	41	LYS	HD3	-0.37	2.75 – 0.45	-8.6
1	A	10	LYS	HG2	-0.81	2.67 – 0.07	-8.4
1	A	10	LYS	HB2	-0.21	3.03 – 0.53	-8.0
1	A	15	PHE	HZ	10.18	9.11 – 4.91	7.5
1	A	10	LYS	HB3	-0.21	3.10 – 0.40	-7.3
1	A	17	HIS	HB2	0.53	4.91 – 1.31	-7.2
1	A	41	LYS	HG2	-0.47	2.67 – 0.07	-7.1
1	A	17	HIS	HB3	0.45	5.00 – 1.10	-6.7
1	A	66	HIS	HB2	0.82	4.91 – 1.31	-6.4
1	A	45	HIS	HB3	0.62	5.00 – 1.10	-6.2
1	A	66	HIS	HB3	0.78	5.00 – 1.10	-5.8
1	A	13	VAL	HG23	-0.81	2.20 – -0.60	-5.8
1	A	13	VAL	HG21	-0.81	2.20 – -0.60	-5.8
1	A	13	VAL	HG22	-0.81	2.20 – -0.60	-5.8
1	A	20	HIS	HB2	1.06	4.91 – 1.31	-5.7
1	A	34	PRO	HD2	1.62	5.45 – 1.85	-5.6
1	A	30	HIS	HB3	0.97	5.00 – 1.10	-5.3
1	A	34	PRO	HD3	1.62	5.52 – 1.72	-5.3
1	A	37	ILE	HD13	-0.84	2.13 – -0.77	-5.2
1	A	37	ILE	HD12	-0.84	2.13 – -0.77	-5.2
1	A	37	ILE	HD11	-0.84	2.13 – -0.77	-5.2
1	A	20	HIS	HB3	1.06	5.00 – 1.10	-5.1
1	A	34	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:



7.2 Chemical shift list 2

File name: BMRB entry 4743

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	14
Number of shifts mapped to atoms	0
Number of unparsed shifts	0
Number of shifts with mapping errors	14
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. First 5 (of 14) occurrences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
UNMAPPED	1	HEC	HMC	3.31	-1.0	1
UNMAPPED	1	HEC	HHC	9.38	-1.0	1
UNMAPPED	1	HEC	HBC	1.49	-1.0	1
UNMAPPED	1	HEC	HMD	3.44	-1.0	1
UNMAPPED	1	HEC	HHB	8.95	-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 776. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/336 (0%)	0/134 (0%)	0/136 (0%)	0/66 (0%)
Sidechain	0/381 (0%)	0/225 (0%)	0/140 (0%)	0/16 (0%)
Aromatic	0/59 (0%)	0/33 (0%)	0/20 (0%)	0/6 (0%)
Overall	0/776 (0%)	0/392 (0%)	0/296 (0%)	0/88 (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.

7.3 Chemical shift list 3

File name: BMRB entry 4743

Chemical shift list name: *assigned_chem_shift_list_3*

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	18
Number of shifts mapped to atoms	0
Number of unparsed shifts	0
Number of shifts with mapping errors	18
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. First 5 (of 18) occurrences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
UNMAPPED	1	HEC	2HAA	3.31	-1.0	1
UNMAPPED	1	HEC	2HBA	3.79	-1.0	1
UNMAPPED	1	HEC	HMC	4.1	-1.0	1
UNMAPPED	1	HEC	HHC	9.76	-1.0	1
UNMAPPED	1	HEC	1HBA	4.08	-1.0	1

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 0%, i.e. 0 atoms were assigned a chemical shift out of a possible 776. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/336 (0%)	0/134 (0%)	0/136 (0%)	0/66 (0%)
Sidechain	0/381 (0%)	0/225 (0%)	0/140 (0%)	0/16 (0%)
Aromatic	0/59 (0%)	0/33 (0%)	0/20 (0%)	0/6 (0%)
Overall	0/776 (0%)	0/392 (0%)	0/296 (0%)	0/88 (0%)

7.3.4 Statistically unusual chemical shifts ⓘ

There are no statistically unusual chemical shifts.

7.3.5 Random Coil Index (RCI) plots ⓘ

No *random coil index* (RCI) plot could be generated from the current chemical shift list (`assigned_chem_shift_list_3`). RCI is only applicable to proteins.

7.4 Chemical shift list 4

File name: BMRB entry 4743

Chemical shift list name: *assigned_chem_shift_list_4*

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	17
Number of shifts mapped to atoms	0
Number of unparsed shifts	0
Number of shifts with mapping errors	17
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. First 5 (of 17) occurrences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
UNMAPPED	1	HEC	2HAD	4.0	-1.0	1
UNMAPPED	1	HEC	HMC	3.02	-1.0	1
UNMAPPED	1	HEC	HHC	9.09	-1.0	1
UNMAPPED	1	HEC	1HBA	3.07	-1.0	1
UNMAPPED	1	HEC	1HAA	4.54	-1.0	1

7.4.2 Chemical shift referencing ⓘ

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

7.4.3 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 776. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/336 (0%)	0/134 (0%)	0/136 (0%)	0/66 (0%)
Sidechain	0/381 (0%)	0/225 (0%)	0/140 (0%)	0/16 (0%)
Aromatic	0/59 (0%)	0/33 (0%)	0/20 (0%)	0/6 (0%)
Overall	0/776 (0%)	0/392 (0%)	0/296 (0%)	0/88 (0%)

7.4.4 Statistically unusual chemical shifts [i](#)

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_4). RCI is only applicable to proteins.