



# Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 11:57 PM BST

PDB ID : 2KW0  
Title : Solution structure of N-terminal domain of CcmH from Escherichia.coli  
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Deposited on : 2010-03-30

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

<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  
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

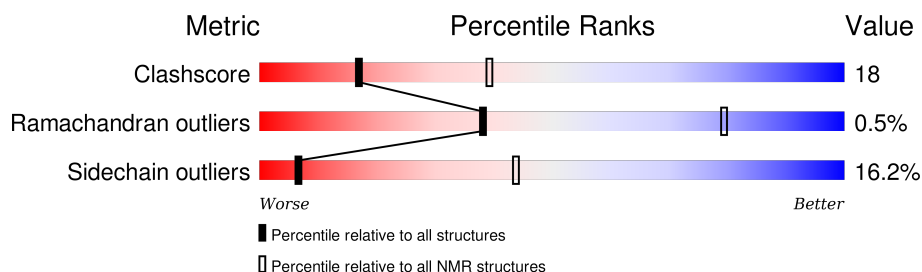
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 83%.

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  $\geq 3$ , 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	90	

## 2 Ensemble composition and analysis ⓘ

This entry contains 20 models. Model 16 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:22-A:95 (74)	0.44	16

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 7 clusters and 2 single-model clusters were found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1308 atoms, of which 648 are hydrogens and 0 are deuteriums.

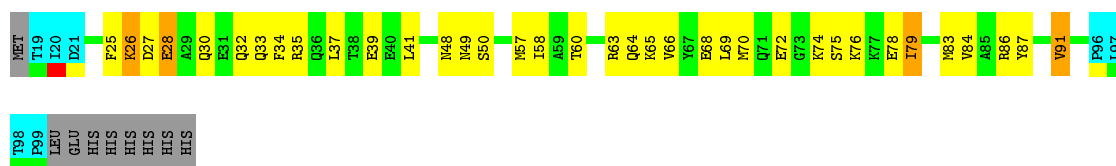
- Molecule 1 is a protein called CcmH protein.

Mol	Chain	Residues	Atoms						Trace
1	A	81	Total	C	H	N	O	S	0
			1308	412	648	111	132	5	

There are 9 discrepancies between the modelled and reference sequences:

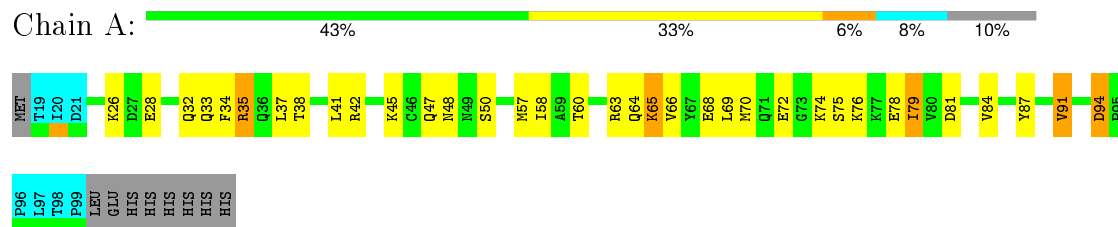
Chain	Residue	Modelled	Actual	Comment	Reference
A	18	MET	-	INITIATING METHIONINE	UNP C5W6T4
A	100	LEU	-	EXPRESSION TAG	UNP C5W6T4
A	101	GLU	-	EXPRESSION TAG	UNP C5W6T4
A	102	HIS	-	EXPRESSION TAG	UNP C5W6T4
A	103	HIS	-	EXPRESSION TAG	UNP C5W6T4
A	104	HIS	-	EXPRESSION TAG	UNP C5W6T4
A	105	HIS	-	EXPRESSION TAG	UNP C5W6T4
A	106	HIS	-	EXPRESSION TAG	UNP C5W6T4
A	107	HIS	-	EXPRESSION TAG	UNP C5W6T4





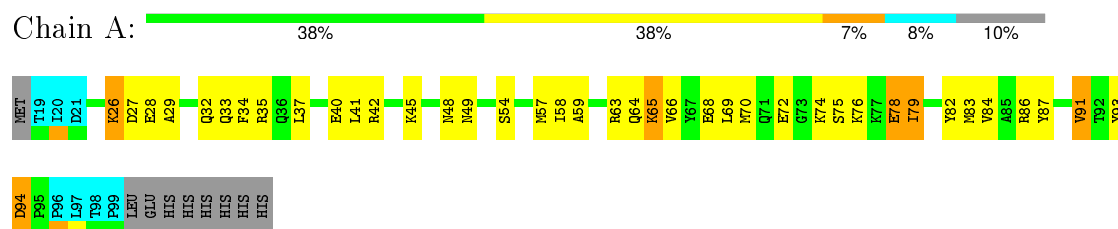
#### 4.2.3 Score per residue for model 3

- Molecule 1: CcmH protein



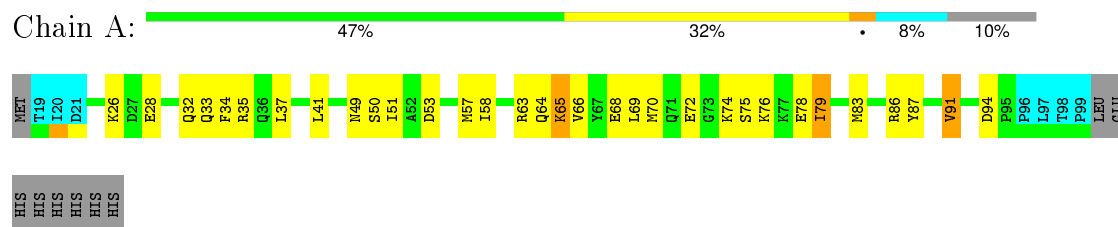
#### 4.2.4 Score per residue for model 4

- Molecule 1: CcmH protein



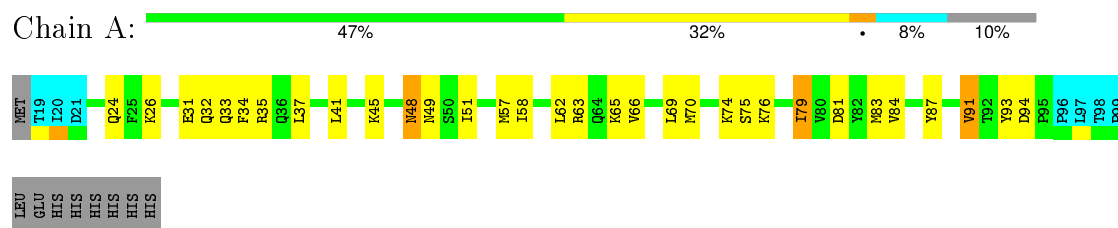
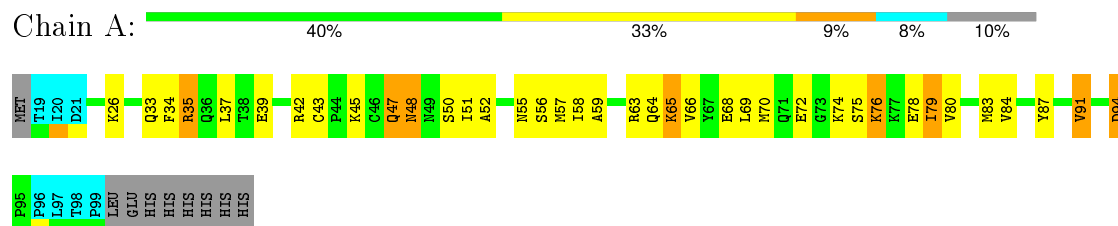
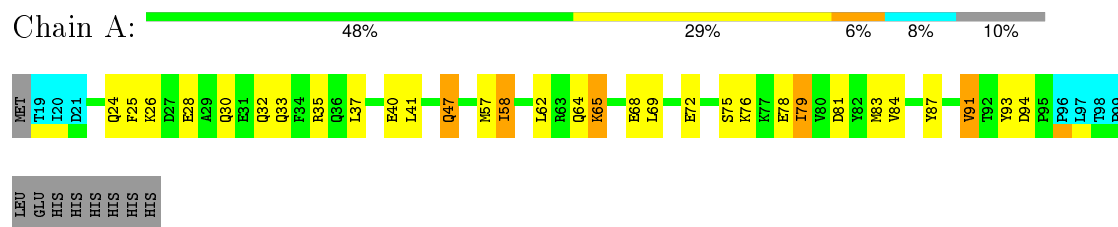
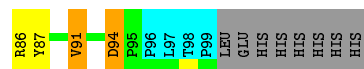
#### 4.2.5 Score per residue for model 5

- Molecule 1: CcmH protein



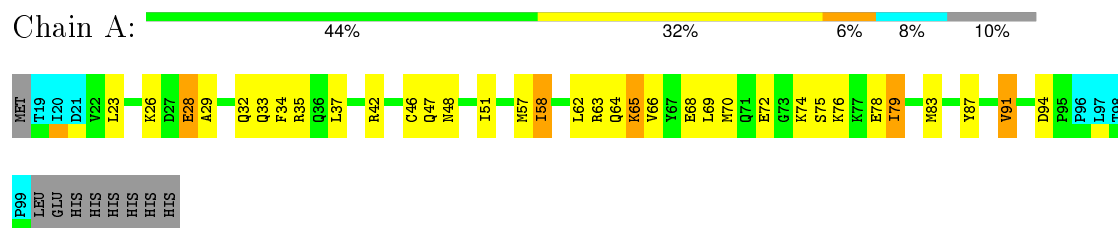
#### 4.2.6 Score per residue for model 6

- Molecule 1: CcmH protein



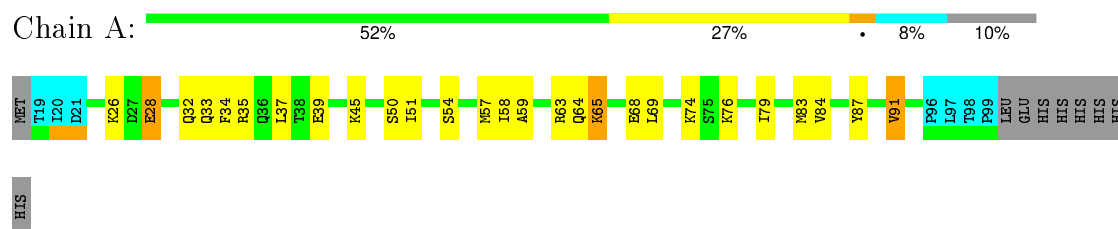
### 4.2.10 Score per residue for model 10

- Molecule 1: CcmH protein



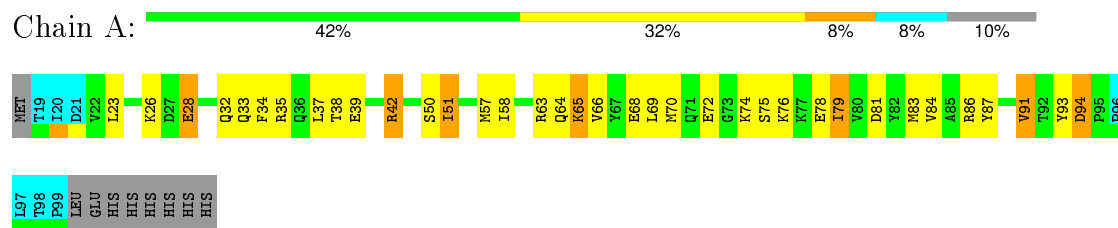
### 4.2.11 Score per residue for model 11

- Molecule 1: CcmH protein



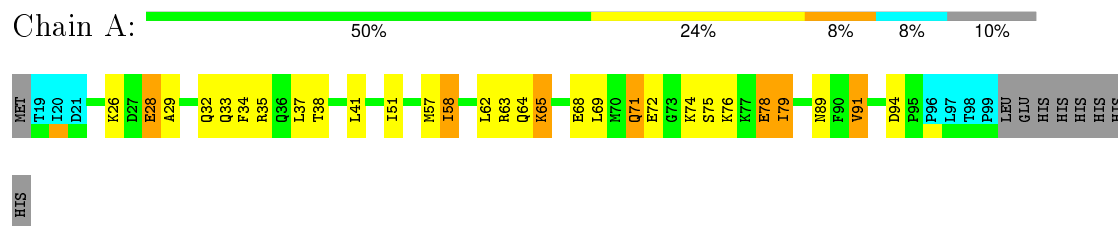
### 4.2.12 Score per residue for model 12

- Molecule 1: CcmH protein



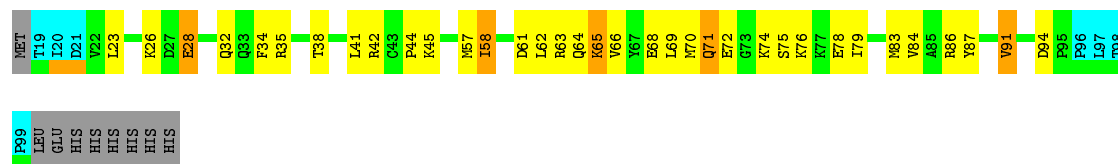
### 4.2.13 Score per residue for model 13

- Molecule 1: CcmH protein









## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing*.

Of the 200 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
CNS	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)	BMRB entry 16802
Number of chemical shift lists	1
Total number of shifts	962
Number of shifts mapped to atoms	962
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	83%

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](#)

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	608	594	594	22±5
All	All	12160	11880	11880	436

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 18.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:41:LEU:HD11	1:A:91:VAL:HB	0.86	1.48	5	13
1:A:69:LEU:HD22	1:A:74:LYS:HG3	0.71	1.62	18	19
1:A:74:LYS:HB3	1:A:78:GLU:HG2	0.68	1.65	4	2
1:A:29:ALA:HA	1:A:32:GLN:HE21	0.68	1.49	10	3
1:A:25:PHE:HA	1:A:30:GLN:NE2	0.67	2.05	15	7
1:A:39:GLU:HA	1:A:50:SER:HB3	0.65	1.69	17	5
1:A:45:LYS:HE3	1:A:47:GLN:HB2	0.62	1.70	15	1
1:A:62:LEU:HD13	1:A:91:VAL:HG12	0.62	1.71	6	5
1:A:49:ASN:HB2	1:A:53:ASP:HB2	0.62	1.71	5	1
1:A:47:GLN:HA	1:A:47:GLN:HE21	0.62	1.53	7	2
1:A:45:LYS:HD3	1:A:48:ASN:HB2	0.62	1.70	9	4
1:A:81:ASP:HA	1:A:84:VAL:HG22	0.60	1.74	3	8
1:A:34:PHE:CZ	1:A:63:ARG:HD2	0.59	2.32	8	15
1:A:28:GLU:O	1:A:32:GLN:HG3	0.59	1.98	20	18

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:65:LYS:HE2	1:A:69:LEU:HG	0.59	1.72	5	8
1:A:43:CYS:SG	1:A:50:SER:HA	0.59	2.37	18	1
1:A:35:ARG:O	1:A:39:GLU:HG3	0.59	1.98	15	2
1:A:69:LEU:HB3	1:A:79:ILE:HG13	0.56	1.76	3	13
1:A:62:LEU:HB3	1:A:83:MET:HE1	0.56	1.77	9	2
1:A:75:SER:H	1:A:78:GLU:HB3	0.56	1.61	2	16
1:A:33:GLN:HE21	1:A:37:LEU:HD11	0.56	1.61	6	1
1:A:66:VAL:O	1:A:70:MET:HG3	0.56	2.01	14	16
1:A:59:ALA:O	1:A:63:ARG:HG2	0.56	2.00	4	1
1:A:35:ARG:HH21	1:A:52:ALA:HB1	0.55	1.61	8	1
1:A:34:PHE:HZ	1:A:63:ARG:HD2	0.55	1.62	4	8
1:A:43:CYS:SG	1:A:49:ASN:HA	0.55	2.42	14	1
1:A:68:GLU:O	1:A:72:GLU:HG3	0.54	2.02	5	17
1:A:31:GLU:O	1:A:35:ARG:HD2	0.54	2.03	15	2
1:A:64:GLN:O	1:A:68:GLU:HG3	0.54	2.02	20	17
1:A:79:ILE:O	1:A:83:MET:HG2	0.54	2.02	14	10
1:A:26:LYS:HD3	1:A:27:ASP:HB3	0.53	1.79	4	1
1:A:38:THR:HG21	1:A:63:ARG:HD2	0.53	1.80	15	1
1:A:69:LEU:HD13	1:A:79:ILE:HA	0.53	1.80	1	8
1:A:41:LEU:HD23	1:A:51:ILE:HD13	0.53	1.80	19	1
1:A:87:TYR:O	1:A:91:VAL:HG13	0.52	2.04	4	18
1:A:65:LYS:HG3	1:A:69:LEU:HD11	0.51	1.82	19	3
1:A:58:ILE:O	1:A:62:LEU:HG	0.51	2.05	20	7
1:A:23:LEU:HD11	1:A:64:GLN:HG2	0.51	1.82	6	2
1:A:65:LYS:HG3	1:A:69:LEU:CD1	0.50	2.37	19	4
1:A:69:LEU:HB3	1:A:74:LYS:HG3	0.50	1.82	14	7
1:A:33:GLN:O	1:A:37:LEU:HG	0.50	2.07	1	17
1:A:34:PHE:CZ	1:A:63:ARG:HB3	0.50	2.42	11	2
1:A:84:VAL:HG11	1:A:93:TYR:HB2	0.49	1.83	16	4
1:A:38:THR:O	1:A:50:SER:HB3	0.49	2.07	3	1
1:A:74:LYS:CB	1:A:78:GLU:HG2	0.49	2.36	4	2
1:A:51:ILE:HA	1:A:54:SER:OG	0.49	2.08	14	3
1:A:39:GLU:HA	1:A:50:SER:CB	0.49	2.37	12	2
1:A:84:VAL:HG12	1:A:91:VAL:HG23	0.49	1.84	15	10
1:A:45:LYS:HE2	1:A:47:GLN:HB2	0.49	1.84	3	1
1:A:56:SER:HB3	1:A:59:ALA:CB	0.49	2.38	19	3
1:A:22:VAL:O	1:A:24:GLN:HG2	0.48	2.07	16	1
1:A:51:ILE:O	1:A:63:ARG:HD3	0.48	2.08	12	6
1:A:39:GLU:HG2	1:A:50:SER:HB3	0.48	1.84	14	1
1:A:41:LEU:O	1:A:50:SER:HB3	0.48	2.08	5	2
1:A:84:VAL:HA	1:A:91:VAL:CG2	0.48	2.39	4	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:75:SER:O	1:A:79:ILE:HD12	0.48	2.08	8	18
1:A:40:GLU:HB3	1:A:93:TYR:HE1	0.48	1.69	15	1
1:A:32:GLN:O	1:A:35:ARG:HB2	0.47	2.09	9	4
1:A:42:ARG:HD2	1:A:94:ASP:CB	0.47	2.39	4	4
1:A:25:PHE:HB3	1:A:30:GLN:HB3	0.47	1.85	1	4
1:A:42:ARG:HD3	1:A:94:ASP:CB	0.47	2.39	6	1
1:A:23:LEU:HD21	1:A:64:GLN:HG2	0.47	1.84	12	1
1:A:75:SER:N	1:A:78:GLU:HB3	0.47	2.25	7	10
1:A:72:GLU:HB3	1:A:74:LYS:HE2	0.47	1.85	19	2
1:A:47:GLN:O	1:A:48:ASN:HB2	0.47	2.10	15	2
1:A:44:PRO:O	1:A:45:LYS:HB3	0.46	2.09	20	1
1:A:60:THR:HA	1:A:63:ARG:HE	0.46	1.71	6	2
1:A:66:VAL:HG12	1:A:70:MET:HE3	0.45	1.87	14	1
1:A:68:GLU:O	1:A:71:GLN:HG3	0.45	2.11	13	2
1:A:32:GLN:HA	1:A:35:ARG:HB2	0.45	1.88	14	1
1:A:59:ALA:HA	1:A:62:LEU:HD12	0.45	1.88	19	1
1:A:86:ARG:HD2	1:A:86:ARG:N	0.44	2.28	2	1
1:A:74:LYS:CA	1:A:78:GLU:HG2	0.44	2.41	13	2
1:A:62:LEU:HB3	1:A:83:MET:CE	0.44	2.43	9	2
1:A:38:THR:CG2	1:A:63:ARG:HG2	0.44	2.43	13	3
1:A:38:THR:OG1	1:A:52:ALA:HB2	0.44	2.13	17	1
1:A:40:GLU:HB3	1:A:93:TYR:CE1	0.44	2.47	15	1
1:A:48:ASN:OD1	1:A:54:SER:HA	0.44	2.13	4	1
1:A:41:LEU:HD12	1:A:42:ARG:H	0.44	1.72	14	1
1:A:42:ARG:HG3	1:A:94:ASP:HB3	0.44	1.89	12	1
1:A:65:LYS:HE3	1:A:65:LYS:HA	0.43	1.90	3	5
1:A:60:THR:O	1:A:64:GLN:HG3	0.43	2.13	2	3
1:A:76:LYS:N	1:A:76:LYS:HE3	0.43	2.28	15	2
1:A:42:ARG:HB2	1:A:92:THR:O	0.43	2.13	19	1
1:A:65:LYS:O	1:A:69:LEU:HG	0.43	2.13	12	1
1:A:40:GLU:HG2	1:A:93:TYR:HE1	0.43	1.73	4	2
1:A:69:LEU:CD2	1:A:74:LYS:HG3	0.42	2.44	8	3
1:A:69:LEU:CB	1:A:79:ILE:HG13	0.42	2.44	3	1
1:A:74:LYS:HB2	1:A:79:ILE:HG13	0.42	1.90	13	1
1:A:38:THR:HG22	1:A:63:ARG:HG2	0.42	1.91	13	2
1:A:31:GLU:O	1:A:34:PHE:HB3	0.42	2.15	9	1
1:A:48:ASN:HB3	1:A:54:SER:HA	0.42	1.92	19	1
1:A:47:GLN:CA	1:A:47:GLN:HE21	0.42	2.25	7	1
1:A:74:LYS:HB3	1:A:78:GLU:CG	0.42	2.44	13	1
1:A:59:ALA:O	1:A:63:ARG:HG3	0.41	2.14	17	2
1:A:42:ARG:HD2	1:A:94:ASP:HB2	0.41	1.92	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:65:LYS:HA	1:A:65:LYS:HE3	0.41	1.93	15	2
1:A:80:VAL:O	1:A:84:VAL:HG13	0.41	2.16	8	1
1:A:78:GLU:O	1:A:82:TYR:HB2	0.41	2.16	4	1
1:A:36:GLN:HB3	1:A:40:GLU:OE2	0.41	2.16	14	1
1:A:65:LYS:HE3	1:A:68:GLU:HB2	0.40	1.93	10	1
1:A:26:LYS:HD2	1:A:27:ASP:OD1	0.40	2.17	2	1
1:A:41:LEU:HD13	1:A:80:VAL:HG23	0.40	1.94	14	1
1:A:72:GLU:CB	1:A:74:LYS:HE2	0.40	2.47	19	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	74/90 (82%)	68±2 (92±2%)	5±2 (7±2%)	0±0 (1±1%)	38	79
All	All	1480/1800 (82%)	1369 (92%)	103 (7%)	8 (1%)	38	79

All 6 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	48	ASN	2
1	A	44	PRO	2
1	A	89	ASN	1
1	A	45	LYS	1
1	A	51	ILE	1
1	A	50	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	68/84 (81%)	57±2 (84±2%)	11±2 (16±2%)	7	44
All	All	1360/1680 (81%)	1140 (84%)	220 (16%)	7	44

All 23 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	58	ILE	20
1	A	57	MET	20
1	A	76	LYS	20
1	A	79	ILE	20
1	A	26	LYS	20
1	A	65	LYS	20
1	A	91	VAL	19
1	A	35	ARG	18
1	A	94	ASP	16
1	A	28	GLU	12
1	A	83	MET	7
1	A	86	ARG	5
1	A	42	ARG	5
1	A	71	GLN	3
1	A	78	GLU	2
1	A	49	ASN	2
1	A	47	GLN	2
1	A	61	ASP	2
1	A	55	ASN	2
1	A	48	ASN	2
1	A	89	ASN	1
1	A	53	ASP	1
1	A	43	CYS	1

### 6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 6.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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 83% for the well-defined parts and 84% for the entire structure.

### 7.1 Chemical shift list 1

File name: BMRB entry 16802

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	962
Number of shifts mapped to atoms	962
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 $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	81	$-0.38 \pm 0.17$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}_\beta$	79	$0.09 \pm 0.11$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}'$	63	$-0.21 \pm 0.13$	None needed ( $< 0.5$ ppm)
$^{15}\text{N}$	66	$0.36 \pm 0.37$	None needed ( $< 0.5$ ppm)

#### 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 83%, i.e. 794 atoms were assigned a chemical shift out of a possible 955. 0 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	327/366 (89%)	135/146 (92%)	131/148 (89%)	61/72 (85%)
Sidechain	443/530 (84%)	279/312 (89%)	157/188 (84%)	7/30 (23%)

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	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Aromatic	24/59 (41%)	24/31 (77%)	0/28 (0%)	0/0 (—%)
Overall	794/955 (83%)	438/489 (90%)	288/364 (79%)	68/102 (67%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 84%, i.e. 867 atoms were assigned a chemical shift out of a possible 1034. 0 out of 11 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	357/397 (90%)	147/158 (93%)	144/162 (89%)	66/77 (86%)
Sidechain	486/578 (84%)	305/340 (90%)	174/208 (84%)	7/30 (23%)
Aromatic	24/59 (41%)	24/31 (77%)	0/28 (0%)	0/0 (—%)
Overall	867/1034 (84%)	476/529 (90%)	318/398 (80%)	73/107 (68%)

#### 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:

