



# Full wwPDB NMR Structure Validation Report ⓘ

Dec 24, 2024 – 04:11 PM EST

PDB ID : 2LJ8  
BMRB ID : 17374  
Title : Solution structure of ADF/Cofilin from trypanosoma brucei  
Authors : Dai, K.; Tu, X.; Zhang, J.; Liao, S.  
Deposited on : 2011-09-08

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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
wwPDB-RCI : v\_1n\_11\_5\_13\_A (Berjanski et al., 2005)  
PANAV : Wang et al. (2010)  
wwPDB-ShiftChecker : v1.2  
BMRB Restraints Analysis : v1.2  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.40

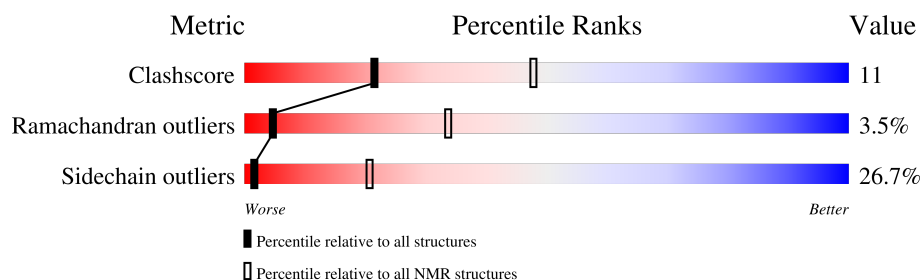
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 81%.

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	210492	14027
Ramachandran outliers	207382	12486
Sidechain outliers	206894	12463

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	144	

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 7 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:14-A:77, A:85-A:144 (124)	0.84	7

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 4 clusters. No single-model clusters were found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2107 atoms, of which 1047 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Cofilin/actin depolymerizing factor, putative.

Mol	Chain	Residues	Atoms						Trace
1	A	136	Total	C	H	N	O	S	0
			2107	662	1047	186	206	6	

There are 8 discrepancies between the modelled and reference sequences:

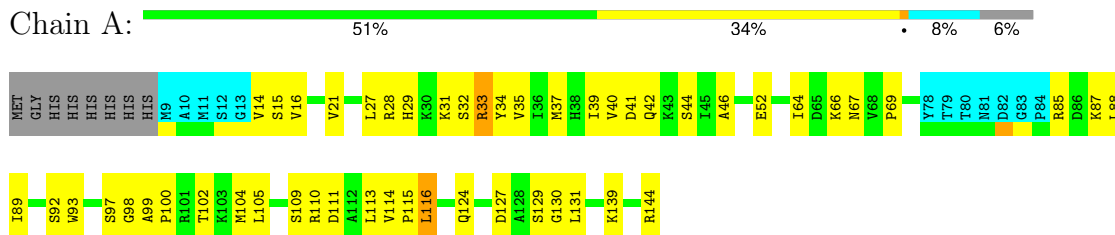
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q580V7
A	2	GLY	-	expression tag	UNP Q580V7
A	3	HIS	-	expression tag	UNP Q580V7
A	4	HIS	-	expression tag	UNP Q580V7
A	5	HIS	-	expression tag	UNP Q580V7
A	6	HIS	-	expression tag	UNP Q580V7
A	7	HIS	-	expression tag	UNP Q580V7
A	8	HIS	-	expression tag	UNP Q580V7

## 4 Residue-property plots

### 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide 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: Cofilin/actin depolymerizing factor, putative

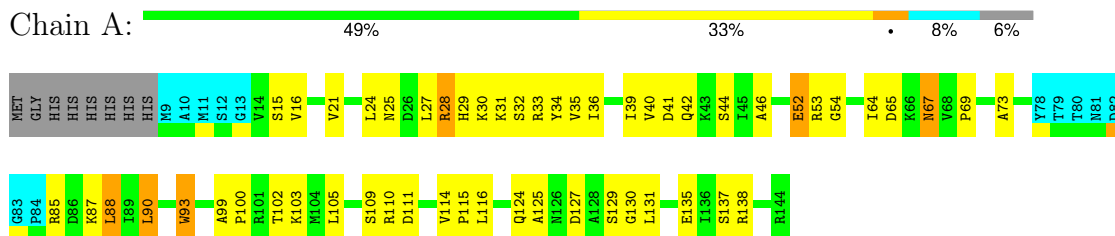


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

#### 4.2.1 Score per residue for model 1

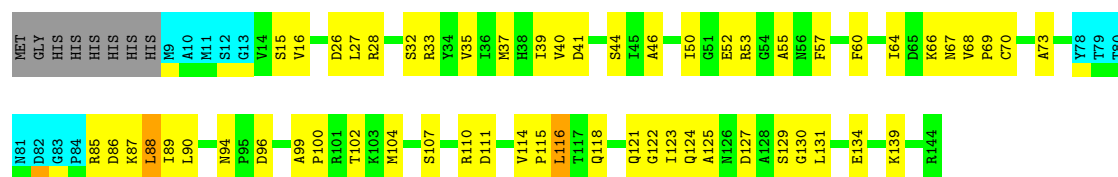
- Molecule 1: Cofilin/actin depolymerizing factor, putative



#### 4.2.2 Score per residue for model 2

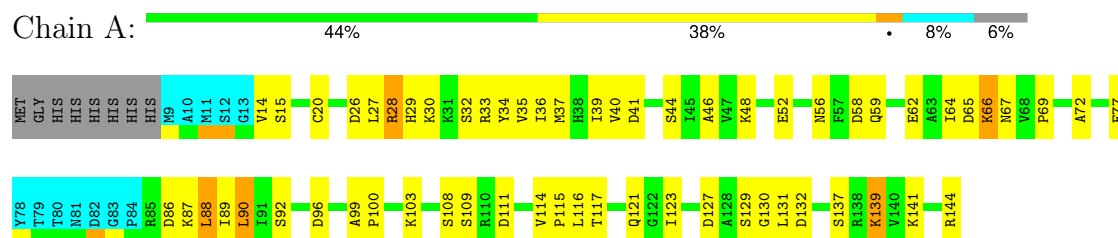
- Molecule 1: Cofilin/actin depolymerizing factor, putative





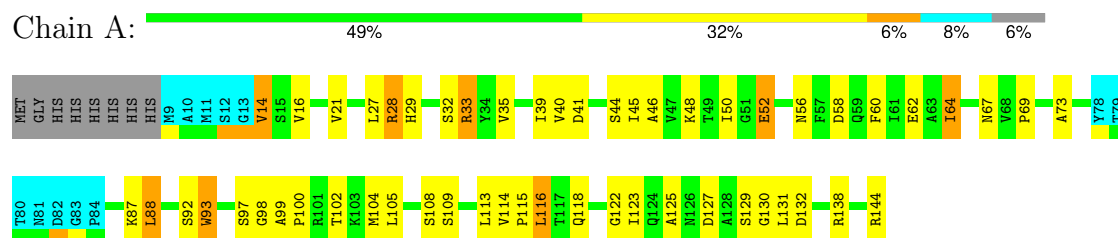
### 4.2.3 Score per residue for model 3

- Molecule 1: Cofilin/actin depolymerizing factor, putative



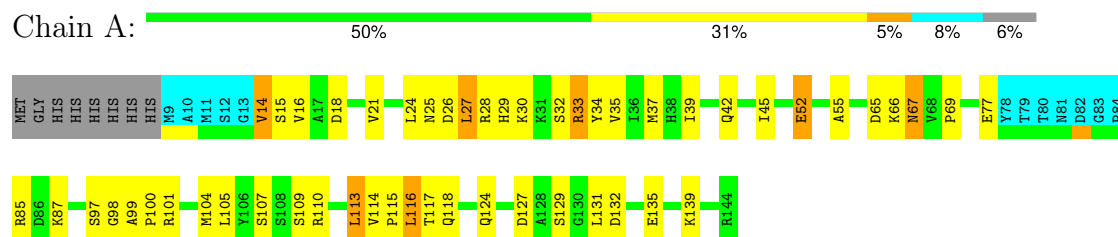
### 4.2.4 Score per residue for model 4

- Molecule 1: Cofilin/actin depolymerizing factor, putative



### 4.2.5 Score per residue for model 5

- Molecule 1: Cofilin/actin depolymerizing factor, putative



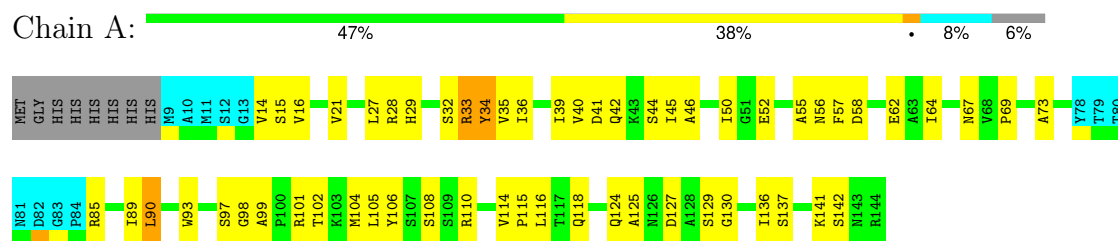
### 4.2.6 Score per residue for model 6

- Molecule 1: Cofilin/actin depolymerizing factor, putative



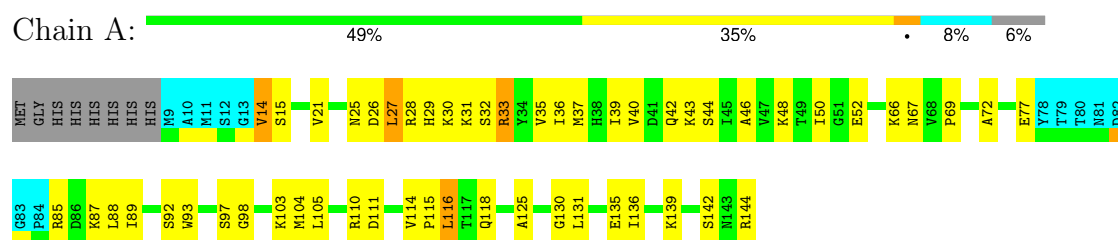
### 4.2.10 Score per residue for model 10

- Molecule 1: Cofilin/actin depolymerizing factor, putative



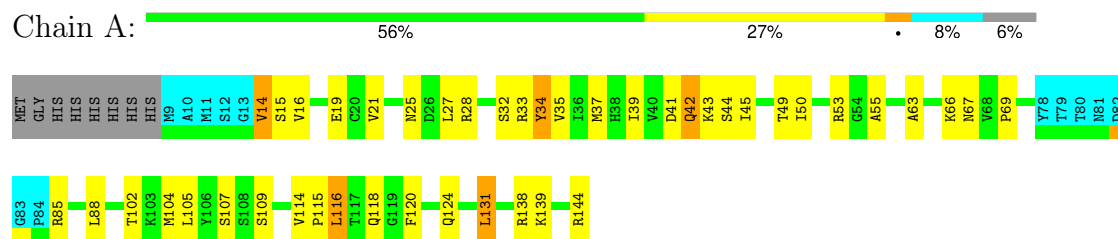
### 4.2.11 Score per residue for model 11

- Molecule 1: Cofilin/actin depolymerizing factor, putative



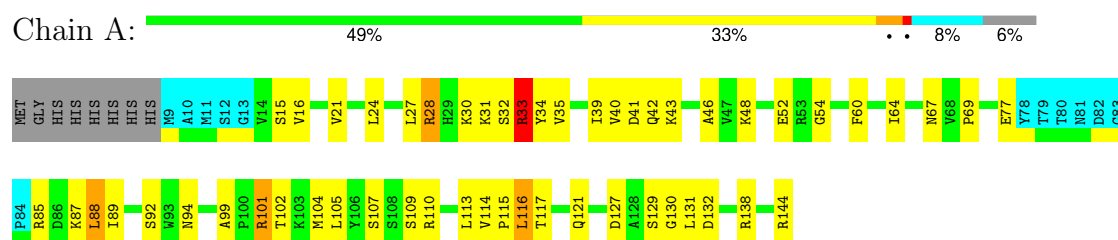
### 4.2.12 Score per residue for model 12

- Molecule 1: Cofilin/actin depolymerizing factor, putative



### 4.2.13 Score per residue for model 13

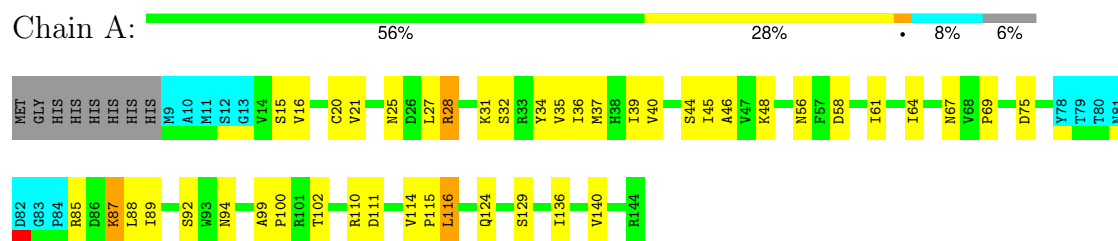
- Molecule 1: Cofilin/actin depolymerizing factor, putative





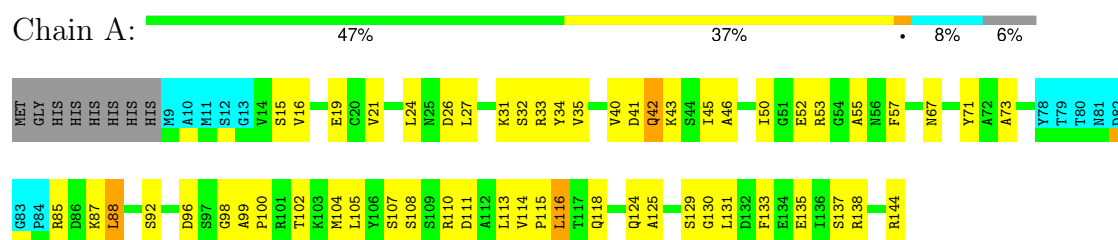
#### 4.2.14 Score per residue for model 14

- Molecule 1: Cofilin/actin depolymerizing factor, putative



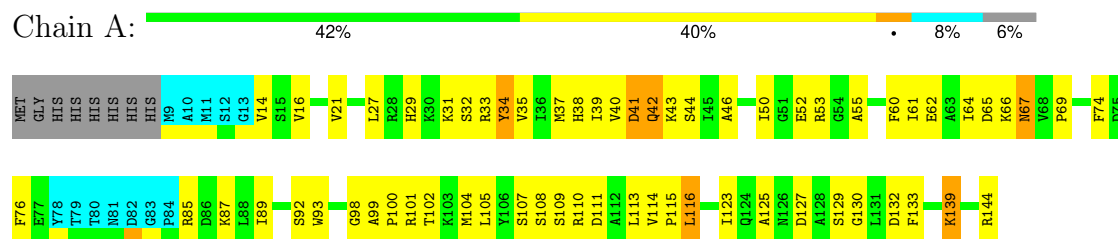
#### 4.2.15 Score per residue for model 15

- Molecule 1: Cofilin/actin depolymerizing factor, putative



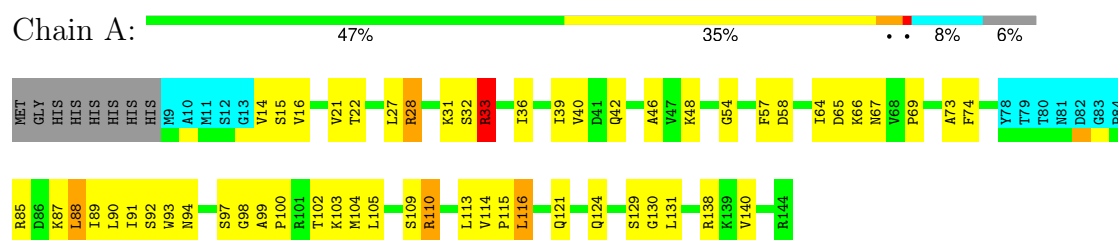
#### 4.2.16 Score per residue for model 16

- Molecule 1: Cofilin/actin depolymerizing factor, putative



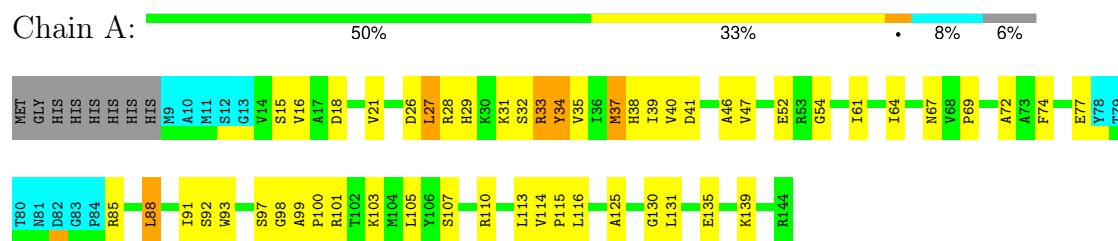
#### 4.2.17 Score per residue for model 17

- Molecule 1: Cofilin/actin depolymerizing factor, putative



### 4.2.18 Score per residue for model 18

- Molecule 1: Cofilin/actin depolymerizing factor, putative



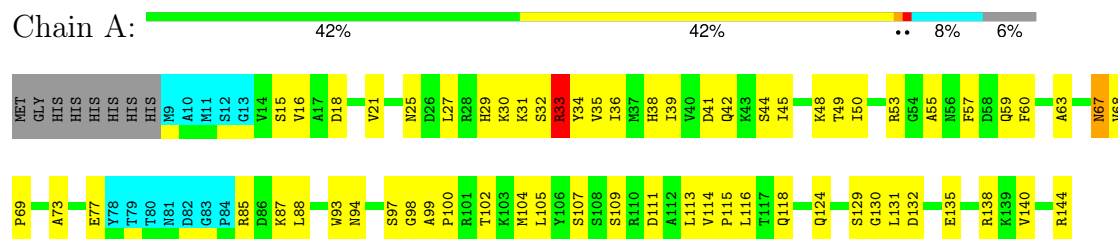
### 4.2.19 Score per residue for model 19

- Molecule 1: Cofilin/actin depolymerizing factor, putative



### 4.2.20 Score per residue for model 20

- Molecule 1: Cofilin/actin depolymerizing factor, putative



## 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
CYANA	structure solution	
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)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	1469
Number of shifts mapped to atoms	1469
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	81%

## 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	976	973	973	22±5
All	All	19520	19460	19460	448

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 11.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:21:VAL:HG13	1:A:116:LEU:HD11	0.94	1.34	6	4
1:A:21:VAL:HG22	1:A:116:LEU:HD11	0.92	1.38	10	13
1:A:16:VAL:HG21	1:A:116:LEU:HD22	0.86	1.48	20	2
1:A:27:LEU:HD13	1:A:35:VAL:HG13	0.82	1.50	14	6
1:A:27:LEU:HD23	1:A:88:LEU:HD11	0.81	1.52	17	1
1:A:27:LEU:HD22	1:A:35:VAL:HG22	0.81	1.51	4	3
1:A:27:LEU:HD23	1:A:88:LEU:HD21	0.81	1.49	2	2
1:A:36:ILE:HG21	1:A:64:ILE:HD11	0.80	1.54	10	4
1:A:36:ILE:HD13	1:A:64:ILE:HD11	0.75	1.58	3	2
1:A:27:LEU:HD13	1:A:35:VAL:HG12	0.70	1.63	6	3
1:A:16:VAL:HG12	1:A:45:ILE:HD12	0.69	1.63	4	1
1:A:27:LEU:HD22	1:A:35:VAL:HG12	0.68	1.65	1	3
1:A:40:VAL:HG21	1:A:46:ALA:HB3	0.64	1.68	14	16
1:A:35:VAL:HG12	1:A:50:ILE:HG23	0.64	1.66	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:93:TRP:CZ3	1:A:131:LEU:HD21	0.64	2.27	11	2
1:A:34:TYR:CE1	1:A:55:ALA:HB3	0.64	2.27	12	3
1:A:99:ALA:HB3	1:A:102:THR:HG23	0.64	1.70	1	4
1:A:14:VAL:HG21	1:A:109:SER:OG	0.64	1.93	16	1
1:A:122:GLY:O	1:A:123:ILE:HD13	0.63	1.93	2	2
1:A:14:VAL:HG21	1:A:109:SER:HB2	0.62	1.72	19	2
1:A:28:ARG:HA	1:A:88:LEU:HD12	0.62	1.69	13	4
1:A:36:ILE:HD11	1:A:60:PHE:CE2	0.62	2.30	20	2
1:A:73:ALA:HB2	1:A:90:LEU:HD12	0.62	1.72	10	2
1:A:35:VAL:HG12	1:A:50:ILE:CG2	0.61	2.24	10	1
1:A:36:ILE:HG21	1:A:64:ILE:CD1	0.60	2.26	1	3
1:A:28:ARG:HB3	1:A:88:LEU:HD12	0.60	1.71	3	2
1:A:45:ILE:HG21	1:A:71:TYR:OH	0.60	1.96	19	2
1:A:27:LEU:HD11	1:A:34:TYR:CA	0.60	2.26	18	3
1:A:39:ILE:HD12	1:A:69:PRO:HG2	0.60	1.72	3	14
1:A:16:VAL:HG21	1:A:116:LEU:CD2	0.60	2.25	20	2
1:A:34:TYR:OH	1:A:55:ALA:HB3	0.59	1.97	7	5
1:A:28:ARG:CB	1:A:88:LEU:HD12	0.59	2.26	3	1
1:A:123:ILE:HD13	1:A:139:LYS:HD2	0.59	1.74	7	1
1:A:55:ALA:HB3	1:A:60:PHE:CE2	0.59	2.33	19	1
1:A:16:VAL:HG11	1:A:116:LEU:HG	0.59	1.75	7	5
1:A:27:LEU:HD21	1:A:34:TYR:HA	0.58	1.74	1	3
1:A:27:LEU:HD11	1:A:34:TYR:HA	0.58	1.73	14	2
1:A:35:VAL:HG12	1:A:50:ILE:HG13	0.58	1.74	4	4
1:A:27:LEU:HD13	1:A:35:VAL:HG23	0.58	1.74	13	4
1:A:94:ASN:OD1	1:A:102:THR:HG21	0.57	1.99	6	1
1:A:24:LEU:HD13	1:A:117:THR:CG2	0.57	2.29	5	4
1:A:42:GLN:CB	1:A:105:LEU:HD11	0.57	2.28	9	3
1:A:93:TRP:CH2	1:A:131:LEU:HD11	0.57	2.34	11	2
1:A:130:GLY:O	1:A:131:LEU:HD23	0.57	1.99	19	9
1:A:73:ALA:HB1	1:A:88:LEU:CD2	0.57	2.29	15	2
1:A:16:VAL:HG11	1:A:116:LEU:CD2	0.57	2.29	13	6
1:A:27:LEU:HD11	1:A:33:ARG:C	0.56	2.20	4	4
1:A:39:ILE:HD12	1:A:69:PRO:CG	0.56	2.30	6	17
1:A:88:LEU:C	1:A:89:ILE:HD13	0.56	2.21	17	6
1:A:21:VAL:HG22	1:A:116:LEU:HD21	0.56	1.78	19	3
1:A:14:VAL:CG2	1:A:105:LEU:HD22	0.56	2.31	12	2
1:A:42:GLN:CG	1:A:105:LEU:HD21	0.55	2.31	10	1
1:A:42:GLN:CG	1:A:105:LEU:HD11	0.55	2.31	17	3
1:A:110:ARG:O	1:A:114:VAL:HG23	0.55	2.01	10	4
1:A:37:MET:CE	1:A:113:LEU:HD21	0.54	2.32	9	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:89:ILE:HG21	1:A:136:ILE:HD12	0.54	1.79	10	2
1:A:28:ARG:CA	1:A:88:LEU:HD12	0.54	2.33	4	3
1:A:14:VAL:HG21	1:A:109:SER:CB	0.54	2.33	19	2
1:A:74:PHE:CE2	1:A:91:ILE:HD13	0.54	2.37	18	1
1:A:40:VAL:CG2	1:A:46:ALA:HB3	0.54	2.33	14	15
1:A:27:LEU:CD2	1:A:88:LEU:HD11	0.54	2.29	17	1
1:A:39:ILE:HD13	1:A:105:LEU:HD22	0.54	1.79	4	1
1:A:99:ALA:HB3	1:A:102:THR:OG1	0.54	2.02	10	4
1:A:14:VAL:HG11	1:A:109:SER:HB2	0.54	1.79	3	1
1:A:93:TRP:CZ2	1:A:131:LEU:HD11	0.54	2.38	20	1
1:A:73:ALA:HB2	1:A:90:LEU:CD1	0.54	2.33	17	1
1:A:60:PHE:O	1:A:64:ILE:HG22	0.54	2.03	16	5
1:A:73:ALA:HB2	1:A:90:LEU:CD2	0.53	2.33	1	1
1:A:114:VAL:N	1:A:115:PRO:HD2	0.53	2.19	7	20
1:A:123:ILE:HD11	1:A:139:LYS:HG2	0.53	1.80	16	2
1:A:16:VAL:CG1	1:A:45:ILE:HD12	0.53	2.33	4	1
1:A:93:TRP:CE2	1:A:131:LEU:HD11	0.52	2.39	20	1
1:A:89:ILE:HG21	1:A:136:ILE:HG23	0.52	1.82	11	1
1:A:90:LEU:CD1	1:A:117:THR:HG21	0.52	2.34	3	1
1:A:28:ARG:CA	1:A:88:LEU:HD21	0.52	2.35	12	4
1:A:21:VAL:CG1	1:A:116:LEU:HD11	0.52	2.32	9	2
1:A:42:GLN:HG2	1:A:105:LEU:HD21	0.52	1.81	10	1
1:A:109:SER:O	1:A:113:LEU:HD12	0.51	2.04	8	2
1:A:27:LEU:HD11	1:A:33:ARG:O	0.51	2.04	17	2
1:A:99:ALA:HB3	1:A:102:THR:HG22	0.51	1.81	16	2
1:A:72:ALA:HB1	1:A:74:PHE:CE1	0.51	2.41	18	1
1:A:99:ALA:HB3	1:A:102:THR:CG2	0.51	2.35	1	3
1:A:93:TRP:CH2	1:A:131:LEU:HD21	0.51	2.41	11	1
1:A:93:TRP:CH2	1:A:131:LEU:HD23	0.50	2.41	18	2
1:A:101:ARG:O	1:A:105:LEU:HD13	0.50	2.06	18	5
1:A:41:ASP:O	1:A:42:GLN:C	0.50	2.49	16	3
1:A:21:VAL:CG2	1:A:116:LEU:HD11	0.50	2.25	10	4
1:A:28:ARG:HA	1:A:88:LEU:HD21	0.50	1.84	9	4
1:A:42:GLN:HB2	1:A:105:LEU:HD21	0.50	1.82	5	3
1:A:21:VAL:HG13	1:A:116:LEU:CD1	0.50	2.33	9	4
1:A:16:VAL:HA	1:A:45:ILE:HD12	0.50	1.83	20	8
1:A:35:VAL:HG22	1:A:50:ILE:HG13	0.50	1.83	9	3
1:A:74:PHE:CZ	1:A:89:ILE:HD13	0.50	2.40	16	1
1:A:69:PRO:N	1:A:102:THR:HG23	0.50	2.22	12	3
1:A:125:ALA:HB3	1:A:130:GLY:HA2	0.49	1.83	4	10
1:A:16:VAL:HG21	1:A:116:LEU:HG	0.49	1.82	14	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:90:LEU:HD13	1:A:117:THR:HG21	0.49	1.83	3	1
1:A:89:ILE:HD13	1:A:136:ILE:CG2	0.49	2.37	6	1
1:A:34:TYR:OH	1:A:72:ALA:HB1	0.49	2.08	3	1
1:A:14:VAL:HG21	1:A:105:LEU:HD22	0.49	1.83	12	1
1:A:87:LYS:HG2	1:A:140:VAL:HG23	0.48	1.85	14	2
1:A:113:LEU:HD23	1:A:116:LEU:HD23	0.48	1.85	20	2
1:A:37:MET:HG2	1:A:47:VAL:HG12	0.48	1.86	18	1
1:A:69:PRO:CD	1:A:102:THR:HG23	0.48	2.38	6	4
1:A:87:LYS:CG	1:A:140:VAL:HG23	0.47	2.39	14	2
1:A:24:LEU:HD12	1:A:25:ASN:N	0.47	2.24	1	1
1:A:16:VAL:HG11	1:A:116:LEU:HD23	0.47	1.87	19	1
1:A:14:VAL:HG12	1:A:45:ILE:HD11	0.47	1.85	5	1
1:A:122:GLY:C	1:A:123:ILE:HD13	0.47	2.30	2	2
1:A:49:THR:OG1	1:A:63:ALA:HB2	0.47	2.09	19	1
1:A:35:VAL:HG12	1:A:50:ILE:CG1	0.47	2.40	12	2
1:A:64:ILE:HD12	1:A:93:TRP:CZ3	0.47	2.45	4	1
1:A:27:LEU:HD21	1:A:73:ALA:O	0.47	2.10	2	3
1:A:27:LEU:CD2	1:A:88:LEU:HD21	0.47	2.31	2	1
1:A:27:LEU:HD23	1:A:88:LEU:CD2	0.46	2.40	6	2
1:A:89:ILE:HD11	1:A:140:VAL:CG2	0.46	2.40	17	1
1:A:60:PHE:CE2	1:A:131:LEU:HD22	0.46	2.46	4	1
1:A:73:ALA:HB1	1:A:88:LEU:HD21	0.46	1.86	8	2
1:A:28:ARG:HA	1:A:88:LEU:HD11	0.45	1.88	14	2
1:A:42:GLN:HB3	1:A:105:LEU:HD21	0.45	1.88	17	1
1:A:42:GLN:HB3	1:A:105:LEU:HD11	0.45	1.87	17	4
1:A:14:VAL:HG22	1:A:14:VAL:O	0.45	2.10	4	1
1:A:123:ILE:HD13	1:A:139:LYS:CD	0.45	2.42	7	1
1:A:27:LEU:HD22	1:A:35:VAL:HB	0.44	1.90	3	1
1:A:24:LEU:HD13	1:A:117:THR:HG22	0.44	1.89	19	1
1:A:116:LEU:O	1:A:116:LEU:HD13	0.44	2.12	19	1
1:A:16:VAL:HG11	1:A:116:LEU:CG	0.44	2.42	2	2
1:A:49:THR:HG21	1:A:63:ALA:HB1	0.44	1.90	12	1
1:A:124:GLN:NE2	1:A:124:GLN:H	0.44	2.10	19	1
1:A:27:LEU:HD22	1:A:35:VAL:HG23	0.44	1.89	9	1
1:A:42:GLN:HE21	1:A:42:GLN:N	0.43	2.12	16	2
1:A:74:PHE:CE1	1:A:91:ILE:HD11	0.43	2.48	17	1
1:A:60:PHE:HE2	1:A:131:LEU:HD22	0.43	1.73	4	1
1:A:42:GLN:HG2	1:A:105:LEU:HD11	0.43	1.89	6	1
1:A:21:VAL:CG2	1:A:116:LEU:HD21	0.43	2.42	7	2
1:A:114:VAL:N	1:A:115:PRO:CD	0.43	2.82	14	20
1:A:89:ILE:HD11	1:A:140:VAL:HG23	0.43	1.91	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:16:VAL:HG11	1:A:116:LEU:HD21	0.42	1.90	13	1
1:A:38:HIS:CE1	1:A:68:VAL:HG21	0.42	2.49	20	1
1:A:14:VAL:HG11	1:A:109:SER:OG	0.42	2.14	5	1
1:A:14:VAL:HG11	1:A:109:SER:CB	0.42	2.44	5	1
1:A:99:ALA:HB1	1:A:100:PRO:HD2	0.42	1.90	5	13
1:A:73:ALA:HB1	1:A:88:LEU:HD22	0.42	1.92	4	1
1:A:109:SER:OG	1:A:113:LEU:HD12	0.42	2.15	5	1
1:A:36:ILE:HD12	1:A:72:ALA:CB	0.41	2.45	11	1
1:A:34:TYR:CZ	1:A:55:ALA:HB3	0.41	2.50	12	1
1:A:55:ALA:HB3	1:A:60:PHE:CZ	0.41	2.50	19	1
1:A:21:VAL:HG22	1:A:116:LEU:CD1	0.41	2.46	18	1
1:A:71:TYR:HB3	1:A:90:LEU:HD11	0.41	1.91	7	1
1:A:49:THR:OG1	1:A:63:ALA:HB1	0.41	2.14	20	1
1:A:50:ILE:HD12	1:A:52:GLU:HG2	0.40	1.93	2	1
1:A:68:VAL:HG23	1:A:70:CYS:SG	0.40	2.57	2	1
1:A:50:ILE:HD11	1:A:52:GLU:OE2	0.40	2.15	11	1
1:A:16:VAL:CA	1:A:45:ILE:HD12	0.40	2.46	20	1
1:A:21:VAL:HG22	1:A:116:LEU:HD13	0.40	1.94	20	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	123/144 (85%)	108±2 (88±2%)	11±2 (9±2%)	4±1 (3±1%)	5	34
All	All	2460/2880 (85%)	2155 (88%)	219 (9%)	86 (3%)	5	34

All 12 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	67	ASN	19
1	A	33	ARG	18
1	A	52	GLU	13
1	A	98	GLY	13

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Mol	Chain	Res	Type	Models (Total)
1	A	14	VAL	7
1	A	54	GLY	4
1	A	53	ARG	3
1	A	131	LEU	3
1	A	43	LYS	3
1	A	55	ALA	1
1	A	66	LYS	1
1	A	32	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	107/123 (87%)	78±4 (73±3%)	29±4 (27±3%)	1	21
All	All	2140/2460 (87%)	1569 (73%)	571 (27%)	1	21

All 80 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	32	SER	19
1	A	15	SER	17
1	A	85	ARG	17
1	A	87	LYS	16
1	A	116	LEU	16
1	A	44	SER	15
1	A	129	SER	15
1	A	104	MET	15
1	A	41	ASP	14
1	A	31	LYS	13
1	A	37	MET	13
1	A	92	SER	13
1	A	29	HIS	12
1	A	110	ARG	12
1	A	111	ASP	12
1	A	127	ASP	12
1	A	33	ARG	12

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Mol	Chain	Res	Type	Models (Total)
1	A	144	ARG	12
1	A	88	LEU	11
1	A	124	GLN	11
1	A	97	SER	11
1	A	138	ARG	10
1	A	66	LYS	10
1	A	113	LEU	10
1	A	28	ARG	9
1	A	107	SER	9
1	A	139	LYS	9
1	A	48	LYS	9
1	A	58	ASP	9
1	A	30	LYS	8
1	A	65	ASP	8
1	A	103	LYS	8
1	A	118	GLN	8
1	A	67	ASN	7
1	A	135	GLU	7
1	A	26	ASP	7
1	A	94	ASN	7
1	A	77	GLU	7
1	A	108	SER	7
1	A	132	ASP	7
1	A	25	ASN	6
1	A	27	LEU	6
1	A	52	GLU	5
1	A	90	LEU	5
1	A	93	TRP	5
1	A	109	SER	5
1	A	137	SER	5
1	A	57	PHE	5
1	A	121	GLN	5
1	A	56	ASN	5
1	A	141	LYS	5
1	A	34	TYR	5
1	A	62	GLU	4
1	A	18	ASP	4
1	A	42	GLN	4
1	A	61	ILE	4
1	A	53	ARG	3
1	A	96	ASP	3
1	A	43	LYS	3

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Mol	Chain	Res	Type	Models (Total)
1	A	120	PHE	3
1	A	75	ASP	3
1	A	19	GLU	3
1	A	133	PHE	3
1	A	86	ASP	2
1	A	134	GLU	2
1	A	20	CYS	2
1	A	59	GLN	2
1	A	64	ILE	2
1	A	76	PHE	2
1	A	89	ILE	2
1	A	101	ARG	2
1	A	131	LEU	2
1	A	142	SER	2
1	A	38	HIS	2
1	A	71	TYR	1
1	A	74	PHE	1
1	A	106	TYR	1
1	A	24	LEU	1
1	A	105	LEU	1
1	A	22	THR	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 oligosaccharides 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 81% for the well-defined parts and 81% for the entire structure.

### 7.1 Chemical shift list 1

File name: working\_cs.cif

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	1469
Number of shifts mapped to atoms	1469
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

#### 7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	133	$0.19 \pm 0.13$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}_\beta$	127	$0.55 \pm 0.14$	Should be applied
$^{13}\text{C}'$	135	$0.12 \pm 0.16$	None needed ( $< 0.5$ ppm)
$^{15}\text{N}$	130	$-0.20 \pm 0.47$	None needed ( $< 0.5$ ppm)

#### 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 81%, i.e. 1364 atoms were assigned a chemical shift out of a possible 1693. 0 out of 17 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	614/618 (99%)	249/250 (100%)	245/248 (99%)	120/120 (100%)
Sidechain	750/962 (78%)	507/622 (82%)	243/296 (82%)	0/44 (0%)

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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	0/113 (0%)	0/56 (0%)	0/54 (0%)	0/3 (0%)
Overall	1364/1693 (81%)	756/928 (81%)	488/598 (82%)	120/167 (72%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 81%, i.e. 1469 atoms were assigned a chemical shift out of a possible 1824. 0 out of 17 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	671/678 (99%)	273/275 (99%)	268/272 (99%)	130/131 (99%)
Sidechain	798/1024 (78%)	539/663 (81%)	259/316 (82%)	0/45 (0%)
Aromatic	0/122 (0%)	0/60 (0%)	0/59 (0%)	0/3 (0%)
Overall	1469/1824 (81%)	812/998 (81%)	527/647 (81%)	130/179 (73%)

#### 7.1.4 Statistically unusual chemical shifts ⓘ

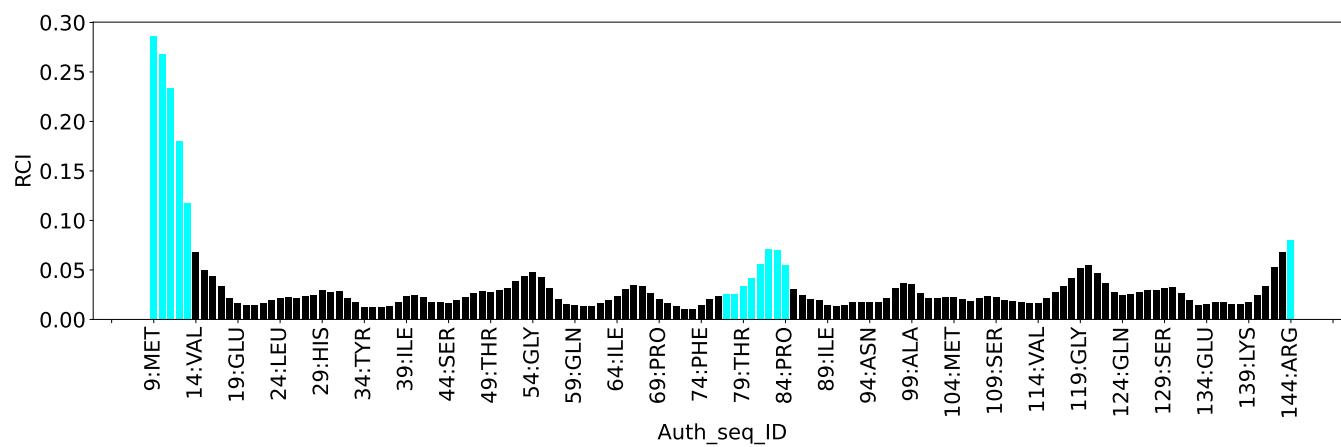
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	51	GLY	HA3	1.38	2.08 – 5.71	-6.9

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

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. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



## 8 NMR restraints analysis

### 8.1 Conformationally restricting restraints

The following table provides the summary of experimentally observed NMR restraints in different categories. Restraints are classified into different categories based on the sequence separation of the atoms involved.

Description	Value
Total distance restraints	1743
Intra-residue ( $ i-j =0$ )	331
Sequential ( $ i-j =1$ )	512
Medium range ( $ i-j >1$ and $ i-j <5$ )	382
Long range ( $ i-j \geq 5$ )	518
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	12.1
Number of long range restraints per residue <sup>1</sup>	3.6

<sup>1</sup>Long range hydrogen bonds and disulfide bonds are counted as long range restraints while calculating the number of long range restraints per residue

### 8.2 Residual restraint violations

This section provides the overview of the restraint violations analysis. The violations are binned as small, medium and large violations based on its absolute value. Average number of violations per model is calculated by dividing the total number of violations in each bin by the size of the ensemble.

#### 8.2.1 Average number of distance violations per model

Distance violations less than 0.1 Å are not included in the calculation.

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	1.4	0.19
0.2-0.5 (Medium)	0.1	0.25
>0.5 (Large)	None	None



### 8.2.2 Average number of dihedral-angle violations per model [i](#)

Dihedral-angle violations less than  $1^\circ$  are not included in the calculation. There are no dihedral-angle violations

## 9 Distance violation analysis ⓘ

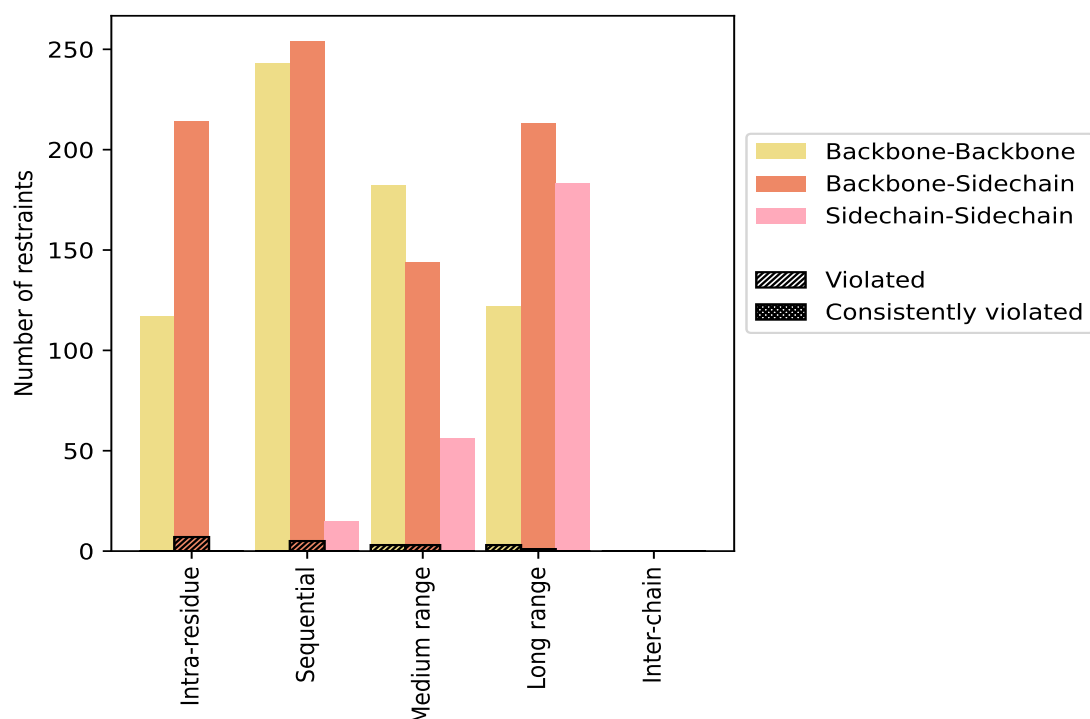
### 9.1 Summary of distance violations ⓘ

The following table shows the summary of distance violations in different restraint categories based on the sequence separation of the atoms involved. Each category is further sub-divided into three sub-categories based on the atoms involved. Violations less than 0.1 Å are not included in the statistics.

Restrains type	Count	% <sup>1</sup>	Violated <sup>3</sup>			Consistently Violated <sup>4</sup>		
			Count	% <sup>2</sup>	% <sup>1</sup>	Count	% <sup>2</sup>	% <sup>1</sup>
<b>Intra-residue (<math> i-j =0</math>)</b>	<b>331</b>	<b>19.0</b>	<b>7</b>	<b>2.1</b>	<b>0.4</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	117	6.7	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	214	12.3	7	3.3	0.4	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
<b>Sequential (<math> i-j =1</math>)</b>	<b>512</b>	<b>29.4</b>	<b>5</b>	<b>1.0</b>	<b>0.3</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	243	13.9	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	254	14.6	5	2.0	0.3	0	0.0	0.0
Sidechain-Sidechain	15	0.9	0	0.0	0.0	0	0.0	0.0
<b>Medium range (<math> i-j &gt;1</math> &amp; <math> i-j &lt;5</math>)</b>	<b>382</b>	<b>21.9</b>	<b>6</b>	<b>1.6</b>	<b>0.3</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	182	10.4	3	1.6	0.2	0	0.0	0.0
Backbone-Sidechain	144	8.3	3	2.1	0.2	0	0.0	0.0
Sidechain-Sidechain	56	3.2	0	0.0	0.0	0	0.0	0.0
<b>Long range (<math> i-j \geq 5</math>)</b>	<b>518</b>	<b>29.7</b>	<b>4</b>	<b>0.8</b>	<b>0.2</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	122	7.0	3	2.5	0.2	0	0.0	0.0
Backbone-Sidechain	213	12.2	1	0.5	0.1	0	0.0	0.0
Sidechain-Sidechain	183	10.5	0	0.0	0.0	0	0.0	0.0
<b>Inter-chain</b>	<b>0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
<b>Hydrogen bond</b>	<b>0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
<b>Disulfide bond</b>	<b>0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
<b>Total</b>	<b>1743</b>	<b>100.0</b>	<b>22</b>	<b>1.3</b>	<b>1.3</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	664	38.1	6	0.9	0.3	0	0.0	0.0
Backbone-Sidechain	825	47.3	16	1.9	0.9	0	0.0	0.0
Sidechain-Sidechain	254	14.6	0	0.0	0.0	0	0.0	0.0

<sup>1</sup> percentage calculated with respect to the total number of distance restraints, <sup>2</sup> percentage calculated with respect to the number of restraints in a particular restraint category, <sup>3</sup> violated in at least one model, <sup>4</sup> violated in all the models

### 9.1.1 Bar chart : Distribution of distance restraints and violations [i](#)



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and disulfied bonds are counted in their appropriate category on the x-axis

## 9.2 Distance violation statistics for each model [i](#)

The following table provides the distance violation statistics for each model in the ensemble. Violations less than 0.1 Å are not included in the statistics.

Model ID	Number of violations						Mean (Å)	Max (Å)	SD <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total				
1	0	0	0	0	0	0	0.0	0.0	0.0	0.0
2	0	0	1	0	0	1	0.12	0.12	0.0	0.12
3	0	0	0	0	0	0	0.0	0.0	0.0	0.0
4	0	0	0	0	0	0	0.0	0.0	0.0	0.0
5	0	0	2	0	0	2	0.12	0.13	0.01	0.12
6	2	0	0	0	0	2	0.16	0.17	0.01	0.16
7	0	0	1	0	0	1	0.11	0.11	0.0	0.11
8	1	0	1	0	0	2	0.18	0.18	0.0	0.18
9	0	0	2	0	0	2	0.12	0.14	0.02	0.12
10	1	2	1	0	0	4	0.15	0.17	0.02	0.16

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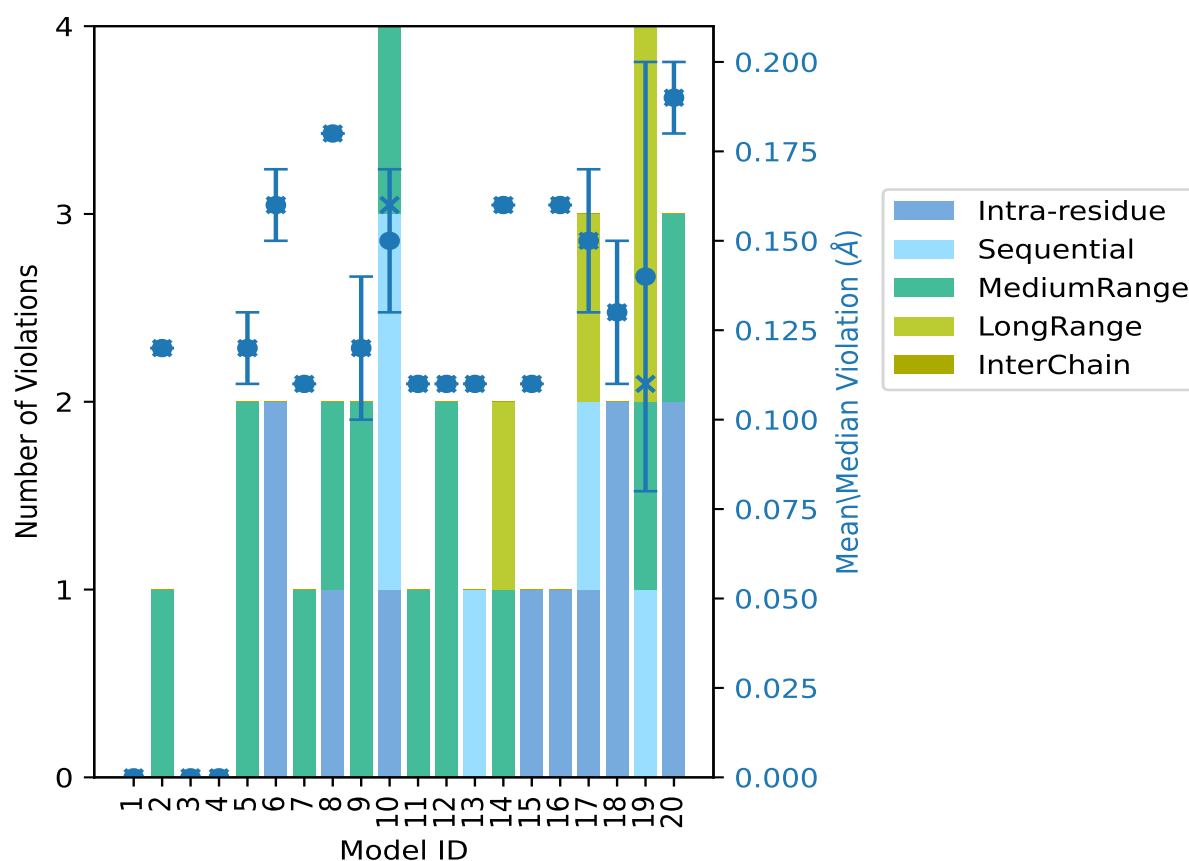
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total				
11	0	0	1	0	0	1	0.11	0.11	0.0	0.11
12	0	0	2	0	0	2	0.11	0.11	0.0	0.11
13	0	1	0	0	0	1	0.11	0.11	0.0	0.11
14	0	0	1	1	0	2	0.16	0.16	0.0	0.16
15	1	0	0	0	0	1	0.11	0.11	0.0	0.11
16	1	0	0	0	0	1	0.16	0.16	0.0	0.16
17	1	1	0	1	0	3	0.15	0.18	0.02	0.15
18	2	0	0	0	0	2	0.13	0.15	0.02	0.13
19	0	1	1	2	0	4	0.14	0.25	0.06	0.11
20	2	0	1	0	0	3	0.19	0.21	0.01	0.19

<sup>1</sup>Intra-residue restraints, <sup>2</sup>Sequential restraints, <sup>3</sup>Medium range restraints, <sup>4</sup>Long range restraints,

<sup>5</sup>Inter-chain restraints, <sup>6</sup>Standard deviation

### 9.2.1 Bar graph : Distance Violation statistics for each model ⓘ



The mean(dot),median(x) and the standard deviation are shown in blue with respect to the y axis on the right

### 9.3 Distance violation statistics for the ensemble

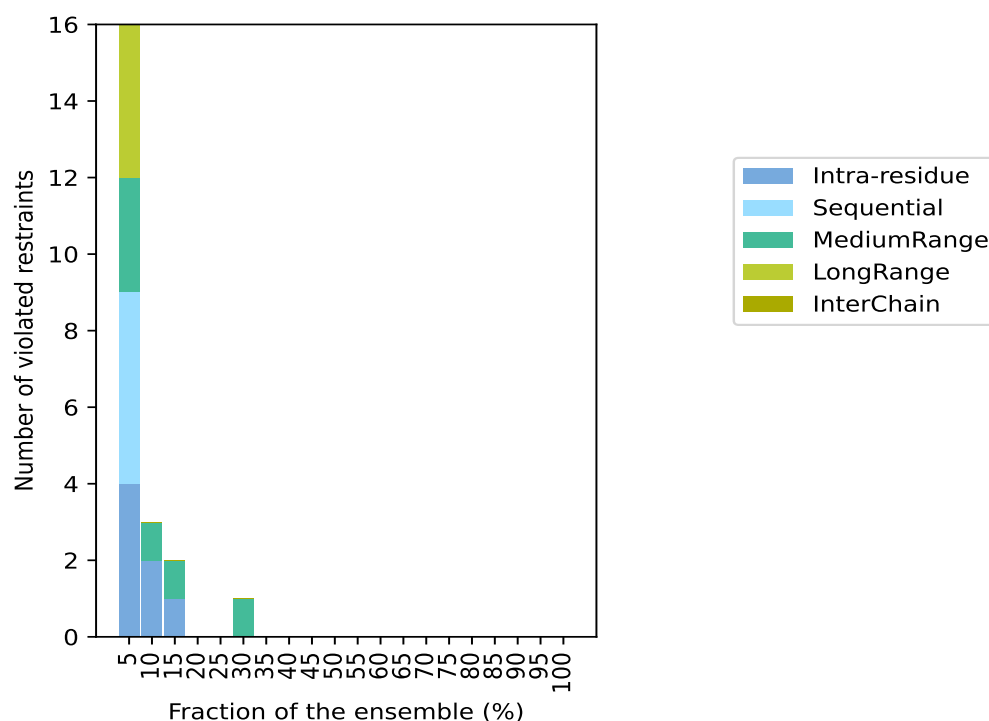
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 1721(IR:324, SQ:507, MR:376, LR:514, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total	Count <sup>6</sup>	%
4	5	3	4	0	16	1	5.0
2	0	1	0	0	3	2	10.0
1	0	1	0	0	2	3	15.0
0	0	0	0	0	0	4	20.0
0	0	0	0	0	0	5	25.0
0	0	1	0	0	1	6	30.0
0	0	0	0	0	0	7	35.0
0	0	0	0	0	0	8	40.0
0	0	0	0	0	0	9	45.0
0	0	0	0	0	0	10	50.0
0	0	0	0	0	0	11	55.0
0	0	0	0	0	0	12	60.0
0	0	0	0	0	0	13	65.0
0	0	0	0	0	0	14	70.0
0	0	0	0	0	0	15	75.0
0	0	0	0	0	0	16	80.0
0	0	0	0	0	0	17	85.0
0	0	0	0	0	0	18	90.0
0	0	0	0	0	0	19	95.0
0	0	0	0	0	0	20	100.0

<sup>1</sup>Intra-residue restraints, <sup>2</sup>Sequential restraints, <sup>3</sup>Medium range restraints, <sup>4</sup>Long range restraints,

<sup>5</sup>Inter-chain restraints, <sup>6</sup> Number of models with violations

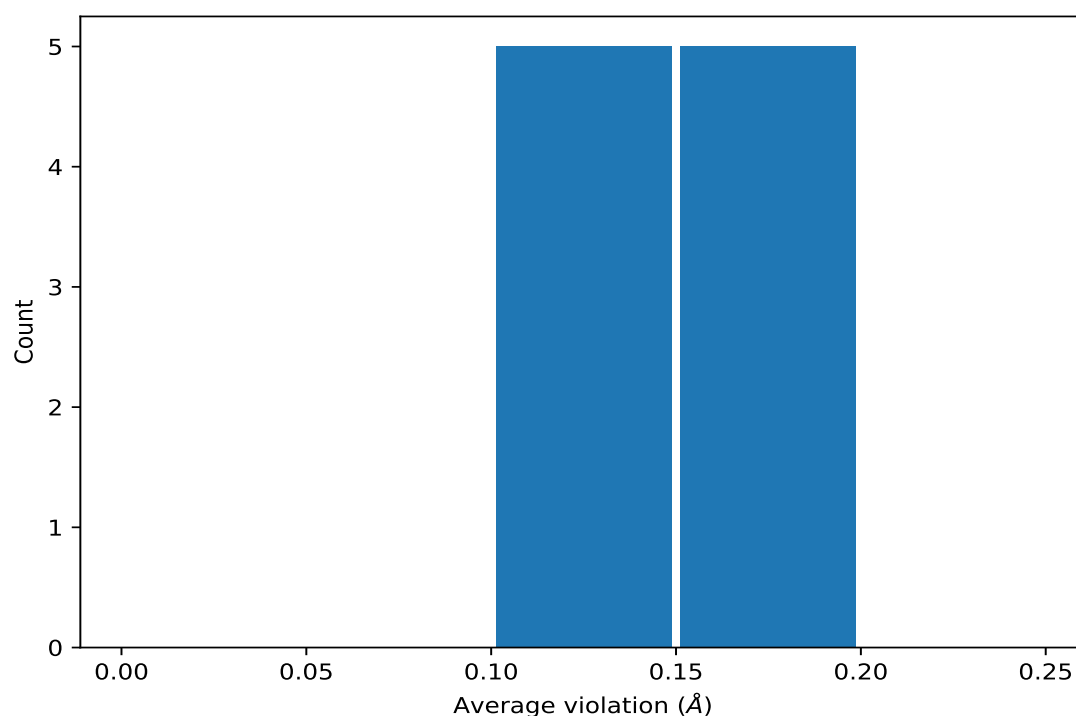
### 9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



## 9.4 Most violated distance restraints in the ensemble [i](#)

### 9.4.1 Histogram : Distribution of mean distance violations [i](#)

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models in the ensemble



#### 9.4.2 Table: Most violated distance restraints [i](#)

The following table provides the mean and the standard deviation of the violation for each restraint sorted by number of violated models and the mean value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

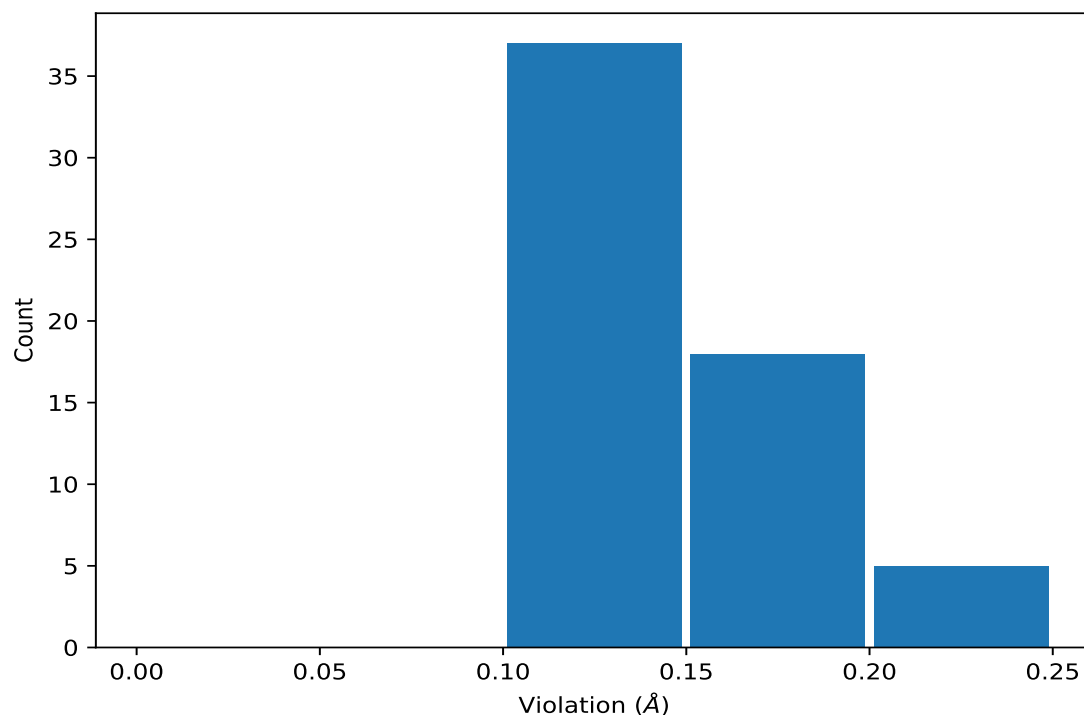
Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(1,38)	1:14:A:VAL:HG21	1:16:A:VAL:HA	6	0.12	0.01	0.11
(1,38)	1:14:A:VAL:HG22	1:16:A:VAL:HA	6	0.12	0.01	0.11
(1,38)	1:14:A:VAL:HG23	1:16:A:VAL:HA	6	0.12	0.01	0.11
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG21	3	0.17	0.01	0.18
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG22	3	0.17	0.01	0.18
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG23	3	0.17	0.01	0.18
(1,903)	1:74:A:PHE:H	1:74:A:PHE:HB2	3	0.17	0.01	0.16
(1,976)	1:82:A:ASP:H	1:82:A:ASP:HB2	2	0.18	0.01	0.18
(1,1122)	1:93:A:TRP:H	1:93:A:TRP:HB3	2	0.13	0.02	0.13
(1,1578)	1:129:A:SER:H	1:125:A:ALA:HA	2	0.1	0.0	0.1

<sup>1</sup>Number of violated models, <sup>2</sup>Standard deviation

## 9.5 All violated distance restraints [i](#)

### 9.5.1 Histogram : Distribution of distance violations [i](#)

The following histogram shows the distribution of the absolute value of the violation for all violated restraints in the ensemble.



### 9.5.2 Table : All distance violations [i](#)

The following table lists the absolute value of the violation for each restraint in the ensemble sorted by its value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,319)	1:32:A:SER:H	1:31:A:LYS:HD2	19	0.25
(1,319)	1:32:A:SER:H	1:31:A:LYS:HD3	19	0.25
(1,958)	1:79:A:THR:H	1:79:A:THR:HG21	20	0.21
(1,958)	1:79:A:THR:H	1:79:A:THR:HG22	20	0.21
(1,958)	1:79:A:THR:H	1:79:A:THR:HG23	20	0.21
(1,976)	1:82:A:ASP:H	1:82:A:ASP:HB2	20	0.19
(1,976)	1:82:A:ASP:H	1:82:A:ASP:HB2	8	0.18
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG21	8	0.18
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG22	8	0.18
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG23	8	0.18

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG21	20	0.18
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG22	20	0.18
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG23	20	0.18
(1,903)	1:74:A:PHE:H	1:74:A:PHE:HB2	17	0.18
(1,1319)	1:111:A:ASP:H	1:110:A:ARG:HB2	10	0.17
(1,268)	1:28:A:ARG:H	1:28:A:ARG:HG2	6	0.17
(1,268)	1:28:A:ARG:H	1:28:A:ARG:HG3	6	0.17
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG21	14	0.16
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG22	14	0.16
(1,972)	1:83:A:GLY:H	1:80:A:THR:HG23	14	0.16
(1,903)	1:74:A:PHE:H	1:74:A:PHE:HB2	10	0.16
(1,903)	1:74:A:PHE:H	1:74:A:PHE:HB2	16	0.16
(1,337)	1:33:A:ARG:H	1:52:A:GLU:H	14	0.16
(1,1662)	1:135:A:GLU:H	1:136:A:ILE:HB	17	0.15
(1,1320)	1:111:A:ASP:H	1:110:A:ARG:HB3	10	0.15
(1,1122)	1:93:A:TRP:H	1:93:A:TRP:HB3	6	0.15
(1,904)	1:74:A:PHE:H	1:74:A:PHE:HB3	18	0.15
(1,38)	1:14:A:VAL:HG21	1:16:A:VAL:HA	9	0.14
(1,38)	1:14:A:VAL:HG22	1:16:A:VAL:HA	9	0.14
(1,38)	1:14:A:VAL:HG23	1:16:A:VAL:HA	9	0.14
(1,964)	1:80:A:THR:H	1:83:A:GLY:H	5	0.13
(1,926)	1:75:A:ASP:H	1:89:A:ILE:HD11	17	0.12
(1,926)	1:75:A:ASP:H	1:89:A:ILE:HD12	17	0.12
(1,926)	1:75:A:ASP:H	1:89:A:ILE:HD13	17	0.12
(1,686)	1:53:A:ARG:H	1:55:A:ALA:H	2	0.12
(1,38)	1:14:A:VAL:HG21	1:16:A:VAL:HA	5	0.12
(1,38)	1:14:A:VAL:HG22	1:16:A:VAL:HA	5	0.12
(1,38)	1:14:A:VAL:HG23	1:16:A:VAL:HA	5	0.12
(1,1122)	1:93:A:TRP:H	1:93:A:TRP:HB3	18	0.11
(1,414)	1:37:A:MET:H	1:71:A:TYR:H	19	0.11
(1,336)	1:33:A:ARG:H	1:52:A:GLU:HA	19	0.11
(1,188)	1:24:A:LEU:H	1:24:A:LEU:HG	15	0.11
(1,38)	1:14:A:VAL:HG21	1:16:A:VAL:HA	7	0.11
(1,38)	1:14:A:VAL:HG22	1:16:A:VAL:HA	7	0.11
(1,38)	1:14:A:VAL:HG23	1:16:A:VAL:HA	7	0.11
(1,38)	1:14:A:VAL:HG21	1:16:A:VAL:HA	10	0.11
(1,38)	1:14:A:VAL:HG22	1:16:A:VAL:HA	10	0.11
(1,38)	1:14:A:VAL:HG23	1:16:A:VAL:HA	10	0.11
(1,38)	1:14:A:VAL:HG21	1:16:A:VAL:HA	11	0.11
(1,38)	1:14:A:VAL:HG22	1:16:A:VAL:HA	11	0.11
(1,38)	1:14:A:VAL:HG23	1:16:A:VAL:HA	11	0.11
(1,38)	1:14:A:VAL:HG21	1:16:A:VAL:HA	12	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,38)	1:14:A:VAL:HG22	1:16:A:VAL:HA	12	0.11
(1,38)	1:14:A:VAL:HG23	1:16:A:VAL:HA	12	0.11
(1,21)	1:13:A:GLY:H	1:14:A:VAL:HG21	13	0.11
(1,21)	1:13:A:GLY:H	1:14:A:VAL:HG22	13	0.11
(1,21)	1:13:A:GLY:H	1:14:A:VAL:HG23	13	0.11
(1,1578)	1:129:A:SER:H	1:125:A:ALA:HA	9	0.1
(1,1578)	1:129:A:SER:H	1:125:A:ALA:HA	12	0.1
(1,233)	1:27:A:LEU:H	1:31:A:LYS:HB3	19	0.1

## 10 Dihedral-angle violation analysis ⓘ

No dihedral-angle restraints found