



# Full wwPDB NMR Structure Validation Report ⓘ

Jun 4, 2023 – 08:34 AM EDT

PDB ID : 2LNE  
BMRB ID : 18162  
Title : Neurotensin 40 structures in water pH 5.5 298 K. NMR data & structures  
Authors : Mukhopadhyay, C.; Khatun, U.  
Deposited on : 2011-12-27

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 : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
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) : 2.33

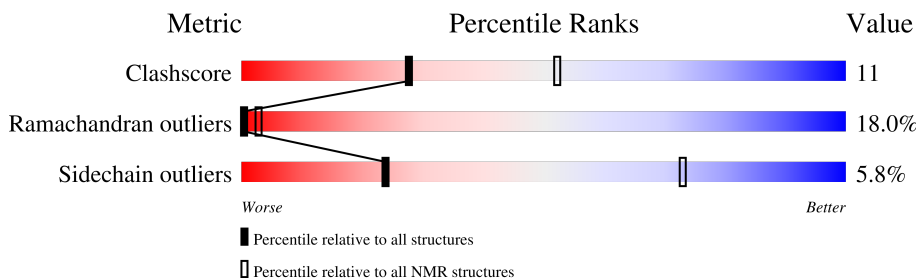
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*


The overall completeness of chemical shifts assignment is 51%.

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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	13	

## 2 Ensemble composition and analysis

This entry contains 40 models.

Cyrange was unable to find well-defined residues.

Error message: Only domains with < 8 residues could be identified.

NmrClust was unable to cluster the ensemble.

Error message: Wrapper check: not enough residues in core to run NmrClust

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 244 atoms, of which 124 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Neurotensin.

Mol	Chain	Residues	Atoms					Trace
			Total	C	H	N	O	
1	A	13	244	78	124	21	21	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLU	GLN	conflict	UNP P30990

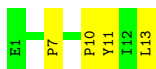
## 4 Residue-property plots

### 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: Neurotensin

Chain A:  69% 31%



### 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: Neurotensin

Chain A:  54% 46%



#### 4.2.2 Score per residue for model 2

- Molecule 1: Neurotensin

Chain A:  54% 46%



### 4.2.3 Score per residue for model 3


- Molecule 1: Neurotensin

Chain A:  46% 54%



### 4.2.4 Score per residue for model 4

- Molecule 1: Neurotensin

Chain A:  77% 23%



### 4.2.5 Score per residue for model 5

- Molecule 1: Neurotensin

Chain A:  62% 38%



### 4.2.6 Score per residue for model 6

- Molecule 1: Neurotensin

Chain A:  69% 15% 15%



### 4.2.7 Score per residue for model 7

- Molecule 1: Neurotensin

Chain A:  69% 31%



#### 4.2.8 Score per residue for model 8

- Molecule 1: Neurotensin

Chain A:  54% 46%



#### 4.2.9 Score per residue for model 9

- Molecule 1: Neurotensin

Chain A:  92% 8%



#### 4.2.10 Score per residue for model 10

- Molecule 1: Neurotensin

Chain A:  62% 38%



#### 4.2.11 Score per residue for model 11


- Molecule 1: Neurotensin

Chain A:  62% 38%



#### 4.2.12 Score per residue for model 12

- Molecule 1: Neurotensin

Chain A:  77% 23%



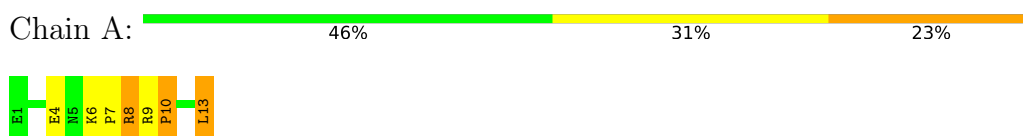
#### 4.2.13 Score per residue for model 13

- Molecule 1: Neurotensin



#### 4.2.14 Score per residue for model 14

- Molecule 1: Neurotensin



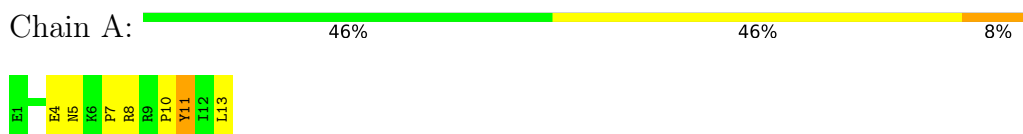
#### 4.2.15 Score per residue for model 15

- Molecule 1: Neurotensin



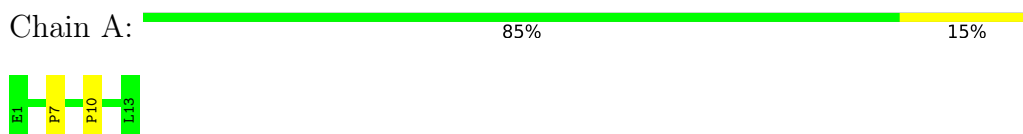
#### 4.2.16 Score per residue for model 16

- Molecule 1: Neurotensin



#### 4.2.17 Score per residue for model 17

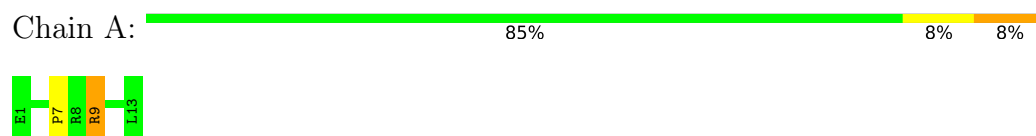
- Molecule 1: Neurotensin





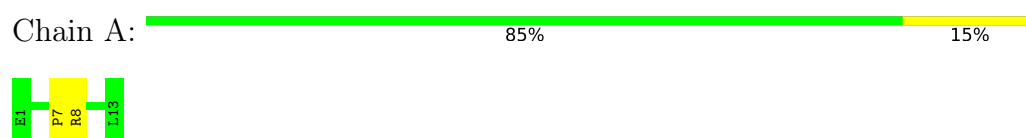
#### 4.2.18 Score per residue for model 18

- Molecule 1: Neurotensin



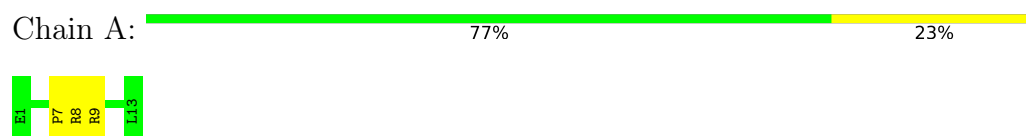
#### 4.2.19 Score per residue for model 19

- Molecule 1: Neurotensin



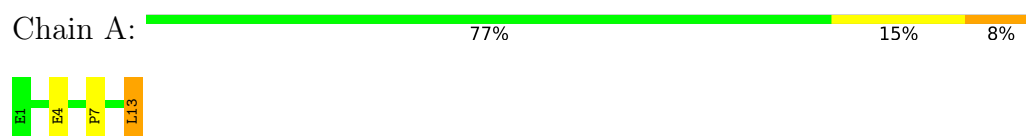
#### 4.2.20 Score per residue for model 20

- Molecule 1: Neurotensin



#### 4.2.21 Score per residue for model 21

- Molecule 1: Neurotensin




#### 4.2.22 Score per residue for model 22

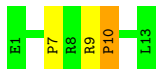
- Molecule 1: Neurotensin



#### 4.2.23 Score per residue for model 23

- Molecule 1: Neurotensin

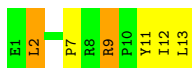
Chain A:  77% 15% 8%



#### 4.2.24 Score per residue for model 24

- Molecule 1: Neurotensin

Chain A:  54% 31% 15%



#### 4.2.25 Score per residue for model 25


- Molecule 1: Neurotensin

Chain A:  69% 23% 8%



#### 4.2.26 Score per residue for model 26


- Molecule 1: Neurotensin

Chain A:  85% 8% 8%



#### 4.2.27 Score per residue for model 27

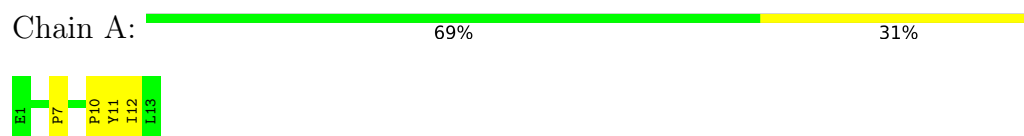
- Molecule 1: Neurotensin

Chain A:  77% 23%



#### 4.2.28 Score per residue for model 28

- Molecule 1: Neurotensin



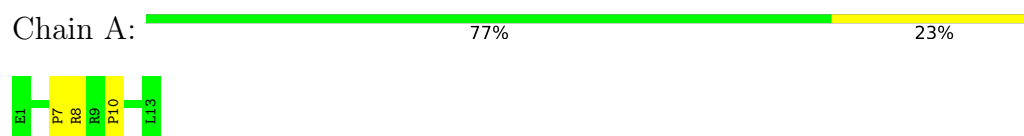
#### 4.2.29 Score per residue for model 29

- Molecule 1: Neurotensin



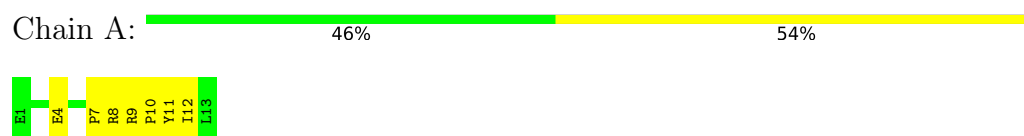
#### 4.2.30 Score per residue for model 30

- Molecule 1: Neurotensin



#### 4.2.31 Score per residue for model 31

- Molecule 1: Neurotensin



#### 4.2.32 Score per residue for model 32

- Molecule 1: Neurotensin



#### 4.2.33 Score per residue for model 33

- Molecule 1: Neurotensin

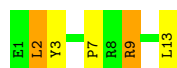
Chain A:  54% 46%



#### 4.2.34 Score per residue for model 34

- Molecule 1: Neurotensin

Chain A:  62% 23% 15%



#### 4.2.35 Score per residue for model 35

- Molecule 1: Neurotensin

Chain A:  69% 31%



#### 4.2.36 Score per residue for model 36

- Molecule 1: Neurotensin

Chain A:  54% 38% 8%



#### 4.2.37 Score per residue for model 37


- Molecule 1: Neurotensin

Chain A:  38% 38% 23%



#### 4.2.38 Score per residue for model 38


- Molecule 1: Neurotensin

Chain A:  77% 15% 8%



#### 4.2.39 Score per residue for model 39


- Molecule 1: Neurotensin

Chain A:  77% 23%



#### 4.2.40 Score per residue for model 40

- Molecule 1: Neurotensin

Chain A:  85% 8% 8%



## 5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 40 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
X-PLOR NIH	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	108
Number of shifts mapped to atoms	108
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	51%

## 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	120	124	124	3±2
All	All	4800	4960	4960	104

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:13:LEU:N	1:A:13:LEU:HD13	0.68	2.04	38	2
1:A:13:LEU:C	1:A:13:LEU:HD22	0.62	2.15	40	1
1:A:12:ILE:O	1:A:12:ILE:HG23	0.58	1.98	6	3
1:A:2:LEU:CD2	1:A:2:LEU:N	0.58	2.66	24	1
1:A:9:ARG:CD	1:A:9:ARG:H	0.57	2.10	34	1
1:A:13:LEU:N	1:A:13:LEU:HD23	0.57	2.15	26	1
1:A:12:ILE:C	1:A:13:LEU:HD13	0.57	2.20	38	1
1:A:10:PRO:O	1:A:12:ILE:N	0.55	2.40	33	4
1:A:13:LEU:CD1	1:A:13:LEU:N	0.55	2.69	40	1
1:A:13:LEU:N	1:A:13:LEU:CD2	0.54	2.70	26	1
1:A:11:TYR:O	1:A:11:TYR:CD2	0.54	2.61	2	6
1:A:4:GLU:H	1:A:4:GLU:CD	0.54	2.06	3	1
1:A:9:ARG:H	1:A:9:ARG:NE	0.54	2.01	18	1
1:A:9:ARG:H	1:A:9:ARG:CD	0.53	2.16	18	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:10:PRO:O	1:A:11:TYR:CB	0.53	2.57	13	1
1:A:4:GLU:O	1:A:5:ASN:ND2	0.53	2.42	22	1
1:A:9:ARG:O	1:A:11:TYR:N	0.52	2.42	2	1
1:A:10:PRO:O	1:A:11:TYR:CG	0.51	2.64	1	1
1:A:4:GLU:OE1	1:A:4:GLU:N	0.51	2.42	27	6
1:A:10:PRO:O	1:A:11:TYR:CD2	0.51	2.64	1	1
1:A:3:TYR:CE2	1:A:4:GLU:O	0.51	2.64	15	1
1:A:13:LEU:HD12	1:A:13:LEU:N	0.50	2.21	13	3
1:A:13:LEU:CD2	1:A:13:LEU:N	0.50	2.74	15	2
1:A:6:LYS:N	1:A:6:LYS:CD	0.49	2.74	10	2
1:A:9:ARG:C	1:A:11:TYR:H	0.49	2.10	2	1
1:A:1:GLU:N	1:A:1:GLU:CD	0.48	2.66	36	4
1:A:10:PRO:C	1:A:11:TYR:CD2	0.48	2.87	1	1
1:A:13:LEU:C	1:A:13:LEU:HD23	0.48	2.29	16	1
1:A:9:ARG:O	1:A:10:PRO:O	0.48	2.31	23	4
1:A:13:LEU:N	1:A:13:LEU:HD22	0.48	2.23	15	2
1:A:11:TYR:CG	1:A:11:TYR:O	0.47	2.66	31	1
1:A:9:ARG:H	1:A:9:ARG:HE	0.46	1.52	18	1
1:A:12:ILE:HG23	1:A:13:LEU:N	0.46	2.26	33	5
1:A:13:LEU:HD23	1:A:13:LEU:H	0.45	1.68	26	1
1:A:10:PRO:C	1:A:12:ILE:N	0.45	2.69	11	3
1:A:9:ARG:O	1:A:11:TYR:CD2	0.45	2.70	24	1
1:A:9:ARG:CD	1:A:9:ARG:N	0.45	2.80	34	1
1:A:5:ASN:OD1	1:A:5:ASN:O	0.44	2.35	1	4
1:A:2:LEU:N	1:A:2:LEU:HD22	0.44	2.26	24	1
1:A:9:ARG:HE	1:A:9:ARG:H	0.44	1.56	11	1
1:A:9:ARG:N	1:A:9:ARG:CD	0.44	2.80	31	1
1:A:12:ILE:CG2	1:A:13:LEU:N	0.44	2.80	33	4
1:A:4:GLU:O	1:A:5:ASN:OD1	0.44	2.36	16	1
1:A:13:LEU:C	1:A:13:LEU:CD2	0.43	2.86	40	1
1:A:8:ARG:CD	1:A:9:ARG:H	0.43	2.25	14	1
1:A:2:LEU:H	1:A:2:LEU:HD23	0.43	1.73	34	1
1:A:6:LYS:H	1:A:6:LYS:CD	0.43	2.24	6	1
1:A:13:LEU:HD13	1:A:13:LEU:H	0.43	1.72	40	1
1:A:11:TYR:O	1:A:11:TYR:CG	0.43	2.72	36	3
1:A:13:LEU:N	1:A:13:LEU:CD1	0.43	2.82	33	2
1:A:13:LEU:H	1:A:13:LEU:CD2	0.42	2.27	21	1
1:A:13:LEU:HD12	1:A:13:LEU:H	0.42	1.75	33	1
1:A:10:PRO:C	1:A:12:ILE:H	0.42	2.18	28	3
1:A:10:PRO:O	1:A:11:TYR:C	0.41	2.57	8	3
1:A:1:GLU:CD	1:A:1:GLU:H1	0.41	2.19	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:13:LEU:C	1:A:13:LEU:HD12	0.41	2.36	14	1
1:A:11:TYR:CD1	1:A:11:TYR:N	0.40	2.88	12	1
1:A:4:GLU:C	1:A:4:GLU:OE1	0.40	2.59	37	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	11/13 (85%)	7±1 (67±9%)	2±1 (15±9%)	2±1 (18±7%)	<b>0</b> <b>3</b>
All	All	440/520 (85%)	296 (67%)	65 (15%)	79 (18%)	<b>0</b> <b>3</b>

All 8 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	7	PRO	40
1	A	8	ARG	13
1	A	11	TYR	9
1	A	10	PRO	8
1	A	12	ILE	3
1	A	4	GLU	3
1	A	3	TYR	2
1	A	6	LYS	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	13/13 (100%)	12±1 (94±6%)	1±1 (6±6%)	<b>24</b> <b>73</b>

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	520/520 (100%)	490 (94%)	30 (6%)	24 73

All 7 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	13	LEU	10
1	A	6	LYS	8
1	A	9	ARG	6
1	A	2	LEU	3
1	A	8	ARG	1
1	A	1	GLU	1
1	A	4	GLU	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 monosaccharides 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 51% for the well-defined parts and 51% 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	108
Number of shifts mapped to atoms	108
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

#### 7.1.2 Chemical shift referencing

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

#### 7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 51%, i.e. 108 atoms were assigned a chemical shift out of a possible 212. 0 out of 2 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	23/61 (38%)	23/24 (96%)	0/26 (0%)	0/11 (0%)
Sidechain	77/133 (58%)	77/85 (91%)	0/40 (0%)	0/8 (0%)
Aromatic	8/18 (44%)	8/8 (100%)	0/10 (0%)	0/0 (—%)
Overall	108/212 (51%)	108/117 (92%)	0/76 (0%)	0/19 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 51%, i.e. 108 atoms were assigned a chemical shift out of a possible 212. 0 out of 2 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	23/61 (38%)	23/24 (96%)	0/26 (0%)	0/11 (0%)
Sidechain	77/133 (58%)	77/85 (91%)	0/40 (0%)	0/8 (0%)
Aromatic	8/18 (44%)	8/8 (100%)	0/10 (0%)	0/0 (—%)
Overall	108/212 (51%)	108/117 (92%)	0/76 (0%)	0/19 (0%)

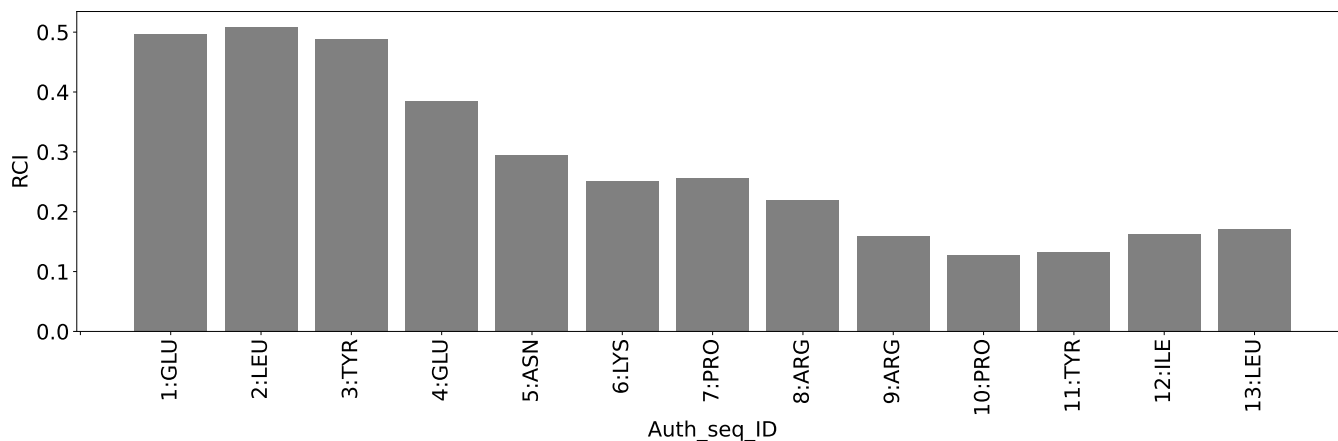
### 7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

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

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	75
Intra-residue ( $ i-j =0$ )	53
Sequential ( $ i-j =1$ )	17
Medium range ( $ i-j >1$ and $ i-j <5$ )	5
Long range ( $ i-j \geq 5$ )	0
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	9
Number of unmapped restraints	0
Number of restraints per residue	6.5
Number of long range restraints per residue <sup>1</sup>	0.0

<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)	1.1	0.2
0.2-0.5 (Medium)	2.9	0.5
>0.5 (Large)	0.0	0.53

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

Dihedral-angle violations less than 1° are not included in the calculation.

Bins (°)	Average number of violations per model	Max (°)
1.0-10.0 (Small)	2.0	9.9
10.0-20.0 (Medium)	0.2	16.0
>20.0 (Large)	1.4	160.5

## 9 Distance violation analysis [i](#)

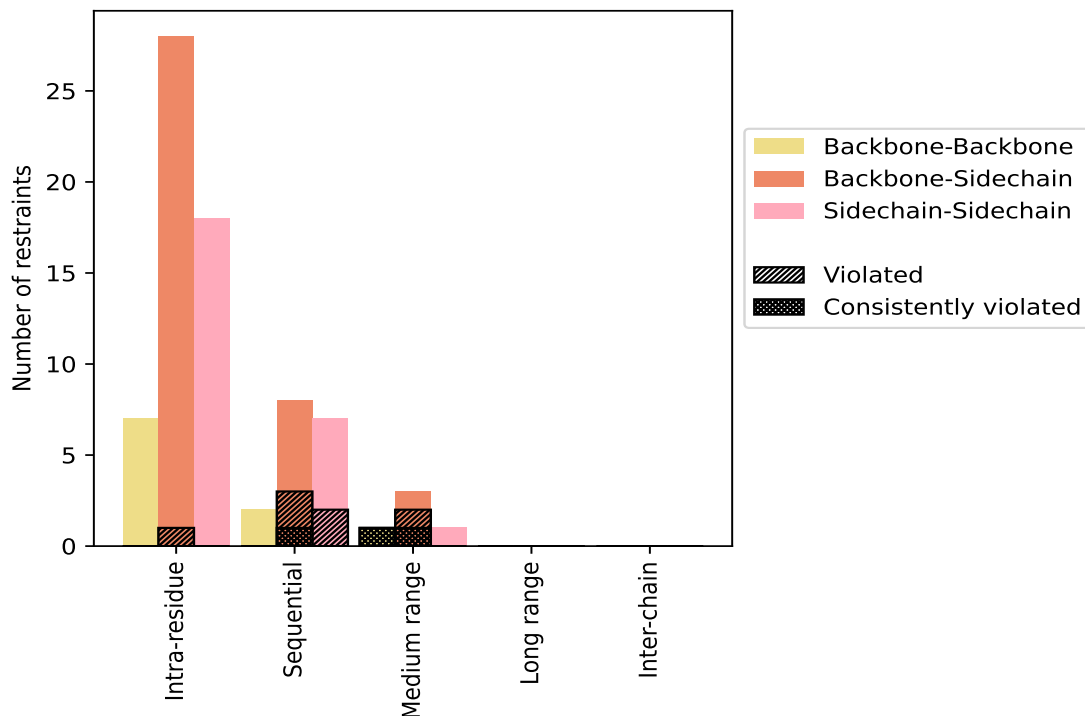
### 9.1 Summary of distance violations [i](#)

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>53</b>	<b>70.7</b>	<b>1</b>	<b>1.9</b>	<b>1.3</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	7	9.3	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	28	37.3	1	3.6	1.3	0	0.0	0.0
Sidechain-Sidechain	18	24.0	0	0.0	0.0	0	0.0	0.0
<b>Sequential (<math> i-j =1</math>)</b>	<b>17</b>	<b>22.7</b>	<b>5</b>	<b>29.4</b>	<b>6.7</b>	<b>1</b>	<b>5.9</b>	<b>1.3</b>
Backbone-Backbone	2	2.7	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	8	10.7	3	37.5	4.0	1	12.5	1.3
Sidechain-Sidechain	7	9.3	2	28.6	2.7	0	0.0	0.0
<b>Medium range (<math> i-j &gt;1</math> &amp; <math> i-j &lt;5</math>)</b>	<b>5</b>	<b>6.7</b>	<b>3</b>	<b>60.0</b>	<b>4.0</b>	<b>2</b>	<b>40.0</b>	<b>2.7</b>
Backbone-Backbone	1	1.3	1	100.0	1.3	1	100.0	1.3
Backbone-Sidechain	3	4.0	2	66.7	2.7	1	33.3	1.3
Sidechain-Sidechain	1	1.3	0	0.0	0.0	0	0.0	0.0
<b>Long range (<math> i-j \geq 5</math>)</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>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>75</b>	<b>100.0</b>	<b>9</b>	<b>12.0</b>	<b>12.0</b>	<b>3</b>	<b>4.0</b>	<b>4.0</b>
Backbone-Backbone	10	13.3	1	10.0	1.3	1	10.0	1.3
Backbone-Sidechain	39	52.0	6	15.4	8.0	2	5.1	2.7
Sidechain-Sidechain	26	34.7	2	7.7	2.7	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 disulfid 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	1	3	0	0	4	0.2	0.25	0.04	0.21
2	0	1	3	0	0	4	0.21	0.28	0.06	0.22
3	0	2	3	0	0	5	0.2	0.27	0.06	0.2
4	0	2	3	0	0	5	0.21	0.29	0.06	0.22
5	1	2	3	0	0	6	0.24	0.5	0.12	0.22
6	0	1	3	0	0	4	0.24	0.39	0.1	0.24
7	0	1	2	0	0	3	0.25	0.28	0.03	0.26
8	0	1	3	0	0	4	0.21	0.29	0.07	0.23
9	1	1	2	0	0	4	0.22	0.3	0.07	0.24
10	0	1	2	0	0	3	0.23	0.26	0.02	0.23
11	0	2	3	0	0	5	0.22	0.29	0.05	0.21

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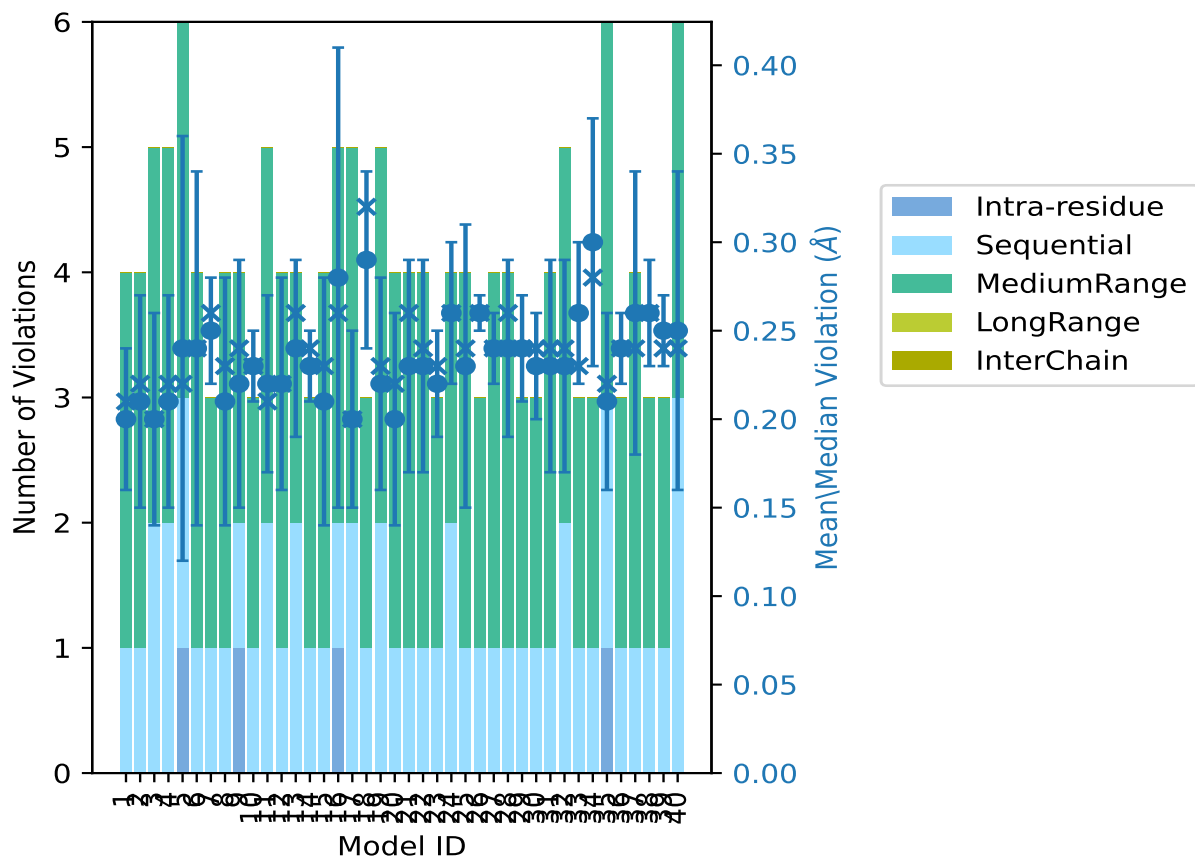
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total				
12	0	1	3	0	0	4	0.22	0.31	0.06	0.22
13	0	2	2	0	0	4	0.24	0.28	0.05	0.26
14	0	1	2	0	0	3	0.23	0.25	0.02	0.24
15	0	1	3	0	0	4	0.21	0.29	0.07	0.23
16	1	1	3	0	0	5	0.28	0.53	0.13	0.26
17	0	2	3	0	0	5	0.2	0.28	0.05	0.2
18	0	1	2	0	0	3	0.29	0.34	0.05	0.32
19	0	2	3	0	0	5	0.22	0.32	0.06	0.23
20	0	1	3	0	0	4	0.2	0.27	0.06	0.22
21	0	1	3	0	0	4	0.23	0.28	0.06	0.26
22	0	1	3	0	0	4	0.23	0.31	0.06	0.24
23	0	1	2	0	0	3	0.22	0.24	0.03	0.23
24	0	2	2	0	0	4	0.26	0.3	0.04	0.26
25	0	1	3	0	0	4	0.23	0.32	0.08	0.24
26	0	1	2	0	0	3	0.26	0.27	0.01	0.26
27	0	1	3	0	0	4	0.24	0.26	0.02	0.24
28	0	1	3	0	0	4	0.24	0.27	0.05	0.26
29	0	1	2	0	0	3	0.24	0.27	0.03	0.24
30	0	1	2	0	0	3	0.23	0.27	0.03	0.24
31	0	1	3	0	0	4	0.23	0.3	0.06	0.24
32	0	2	3	0	0	5	0.23	0.3	0.06	0.24
33	0	1	2	0	0	3	0.26	0.32	0.04	0.23
34	0	1	2	0	0	3	0.3	0.4	0.07	0.28
35	1	2	3	0	0	6	0.21	0.28	0.05	0.22
36	0	1	2	0	0	3	0.24	0.27	0.02	0.24
37	0	1	3	0	0	4	0.26	0.37	0.08	0.24
38	0	1	2	0	0	3	0.26	0.29	0.03	0.26
39	0	1	2	0	0	3	0.25	0.28	0.02	0.24
40	0	3	3	0	0	6	0.25	0.39	0.09	0.24

<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 [i](#)



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 [i](#)

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, 66(IR:52, SQ:12, MR:2, LR:0, 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>	%
0	2	0	0	0	2	1	2.5
0	0	0	0	0	0	2	5.0
0	0	0	0	0	0	3	7.5
1	1	0	0	0	2	4	10.0
0	0	0	0	0	0	5	12.5
0	1	0	0	0	1	6	15.0

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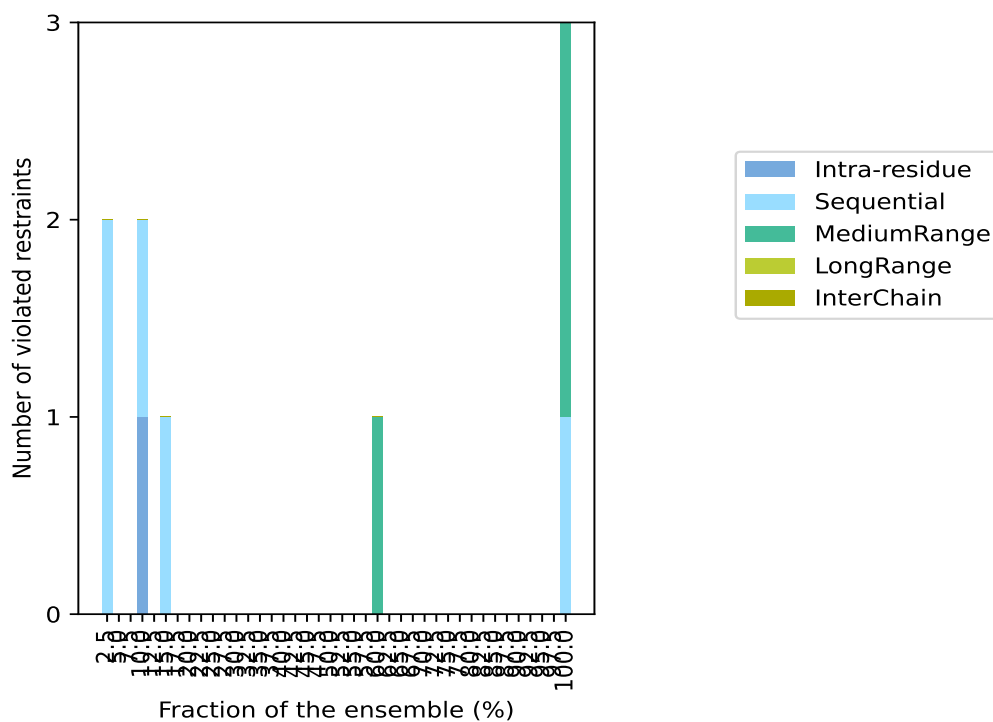
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Number of violated restraints						Fraction of the ensemble	
IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total	Count <sup>6</sup>	%
0	0	0	0	0	0	7	17.5
0	0	0	0	0	0	8	20.0
0	0	0	0	0	0	9	22.5
0	0	0	0	0	0	10	25.0
0	0	0	0	0	0	11	27.5
0	0	0	0	0	0	12	30.0
0	0	0	0	0	0	13	32.5
0	0	0	0	0	0	14	35.0
0	0	0	0	0	0	15	37.5
0	0	0	0	0	0	16	40.0
0	0	0	0	0	0	17	42.5
0	0	0	0	0	0	18	45.0
0	0	0	0	0	0	19	47.5
0	0	0	0	0	0	20	50.0
0	0	0	0	0	0	21	52.5
0	0	0	0	0	0	22	55.0
0	0	0	0	0	0	23	57.5
0	0	1	0	0	1	24	60.0
0	0	0	0	0	0	25	62.5
0	0	0	0	0	0	26	65.0
0	0	0	0	0	0	27	67.5
0	0	0	0	0	0	28	70.0
0	0	0	0	0	0	29	72.5
0	0	0	0	0	0	30	75.0
0	0	0	0	0	0	31	77.5
0	0	0	0	0	0	32	80.0
0	0	0	0	0	0	33	82.5
0	0	0	0	0	0	34	85.0
0	0	0	0	0	0	35	87.5
0	0	0	0	0	0	36	90.0
0	0	0	0	0	0	37	92.5
0	0	0	0	0	0	38	95.0
0	0	0	0	0	0	39	97.5
0	1	2	0	0	3	40	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

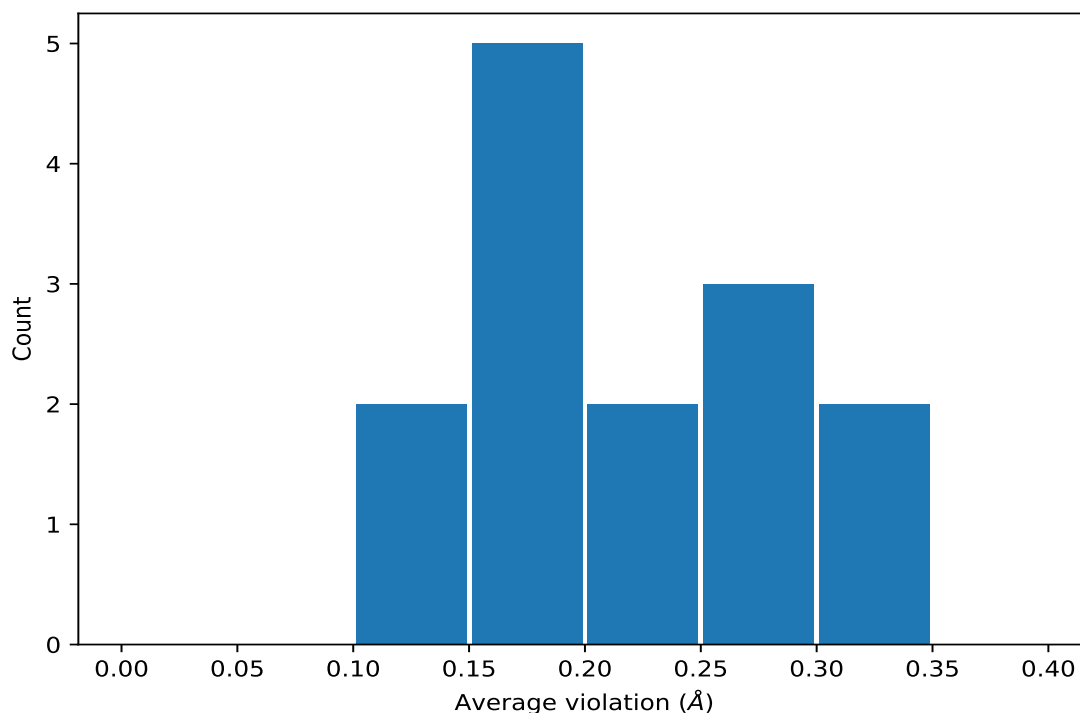
### 9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



## 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.

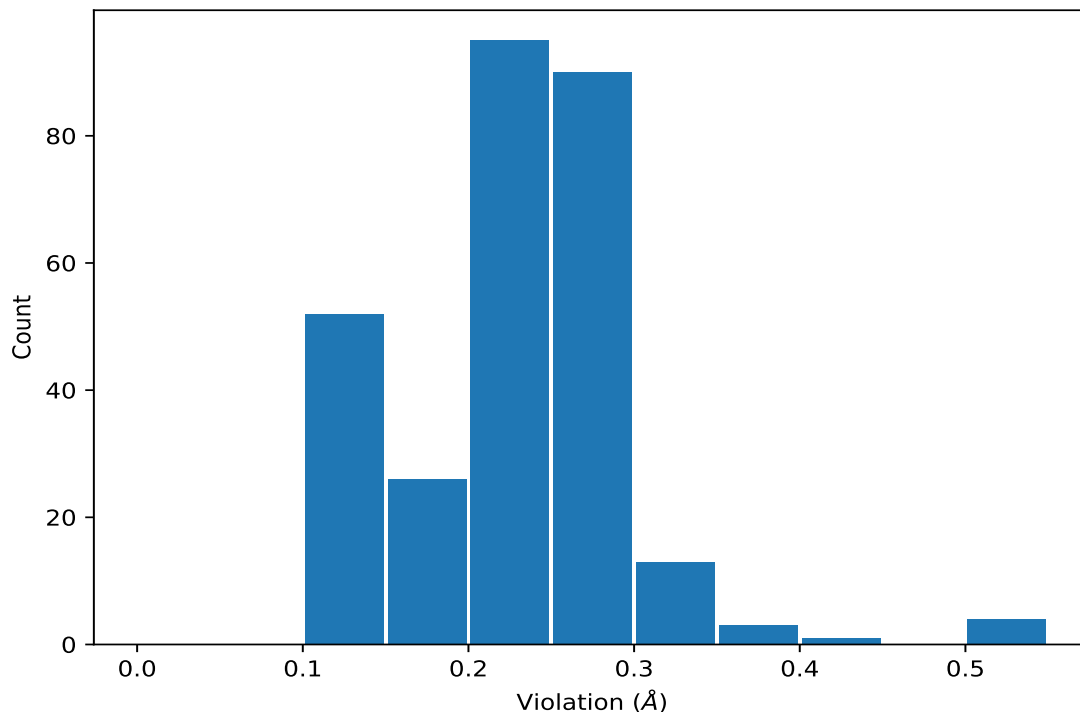
Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	40	0.28	0.04	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	40	0.27	0.03	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	40	0.27	0.03	0.26
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	40	0.22	0.02	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	40	0.22	0.02	0.22
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	24	0.14	0.03	0.13
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	24	0.14	0.03	0.13
(1,54)	1:A:1:GLU:HA	1:A:2:LEU:HG	6	0.16	0.05	0.14
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD2	4	0.34	0.17	0.36
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD3	4	0.34	0.17	0.36
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB2	4	0.16	0.01	0.16
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB3	4	0.16	0.01	0.16
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB2	4	0.16	0.01	0.16
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB3	4	0.16	0.01	0.16

<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)
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD2	16	0.53
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD3	16	0.53
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD2	5	0.5
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD3	5	0.5
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	34	0.4
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	6	0.39
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	40	0.39
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	37	0.37
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	18	0.34
(1,55)	1:A:2:LEU:HA	1:A:1:GLU:HG2	40	0.34

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,55)	1:A:2:LEU:HA	1:A:1:GLU:HG3	40	0.34
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	18	0.32
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	18	0.32
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	25	0.32
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	25	0.32
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	33	0.32
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	33	0.32
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	19	0.32
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	12	0.31
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	12	0.31
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	22	0.31
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	24	0.3
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	24	0.3
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	31	0.3
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	31	0.3
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	9	0.3
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	32	0.3
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	4	0.29
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	4	0.29
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	11	0.29
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	11	0.29
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	16	0.29
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	16	0.29
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	38	0.29
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	38	0.29
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	8	0.29
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	15	0.29
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	24	0.29
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	25	0.29
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	13	0.28
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	13	0.28
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	2	0.28
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	2	0.28
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	7	0.28
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	7	0.28
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	17	0.28
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	17	0.28
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	34	0.28
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	34	0.28
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	13	0.28
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	21	0.28
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	35	0.28

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	39	0.28
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	21	0.27
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	21	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	3	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	3	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	20	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	20	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	28	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	28	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	29	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	29	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	30	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	30	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	36	0.27
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	36	0.27
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	5	0.27
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	26	0.27
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	27	0.26
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	27	0.26
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	28	0.26
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	28	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	9	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	9	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	10	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	10	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	15	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	15	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	26	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	26	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	32	0.26
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	32	0.26
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	3	0.26
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	4	0.26
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	7	0.26
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	16	0.26
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	28	0.26
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	31	0.26
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	38	0.26
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	22	0.25
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	22	0.25
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	26	0.25
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	26	0.25

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	1	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	1	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	6	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	6	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	13	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	13	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	27	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	27	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	35	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	35	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	37	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	37	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	40	0.25
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	40	0.25
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	11	0.25
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	12	0.25
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	14	0.25
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	32	0.24
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	32	0.24
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	37	0.24
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	37	0.24
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	40	0.24
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	40	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	8	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	8	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	14	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	14	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	21	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	21	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	22	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	22	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	39	0.24
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	39	0.24
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	2	0.24
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	17	0.24
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	20	0.24
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	23	0.24
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	27	0.24
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	29	0.24
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	30	0.24
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	36	0.24
(1,54)	1:A:1:GLU:HA	1:A:2:LEU:HG	24	0.24

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	19	0.23
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	19	0.23
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	33	0.23
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	33	0.23
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	34	0.23
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	34	0.23
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	19	0.23
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	19	0.23
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	23	0.23
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	23	0.23
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	10	0.23
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	33	0.23
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD2	35	0.23
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD3	35	0.23
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	4	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	4	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	5	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	5	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	6	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	6	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	8	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	8	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	18	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	18	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	35	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	35	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	36	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	36	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	38	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	38	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	39	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	39	0.22
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB2	5	0.22
(1,63)	1:A:8:ARG:HA	1:A:6:LYS:HB3	5	0.22
(1,62)	1:A:8:ARG:H	1:A:10:PRO:HA	1	0.22
(1,54)	1:A:1:GLU:HA	1:A:2:LEU:HG	32	0.22
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	7	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	7	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	9	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	9	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	10	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	10	0.21

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	11	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	11	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	14	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	14	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	16	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	16	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	24	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	24	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	31	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	31	0.21
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	1	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	1	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	2	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	2	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	3	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	3	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	15	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	15	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	20	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	20	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	25	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	25	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	29	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	29	0.2
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	17	0.2
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	17	0.2
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	27	0.2
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	27	0.2
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	12	0.19
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	12	0.19
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	30	0.19
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	30	0.19
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	17	0.18
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	17	0.18
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG2	23	0.18
(1,65)	1:A:8:ARG:HA	1:A:7:PRO:HG3	23	0.18
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB2	3	0.18
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB3	3	0.18
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB2	3	0.18
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB3	3	0.18
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	11	0.18
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	11	0.18

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	4	0.17
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	4	0.17
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB2	11	0.16
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB3	11	0.16
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB2	11	0.16
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB3	11	0.16
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	19	0.16
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	19	0.16
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	37	0.16
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	37	0.16
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	40	0.16
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	40	0.16
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB2	19	0.15
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB3	19	0.15
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB2	19	0.15
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB3	19	0.15
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB2	35	0.15
(1,64)	1:A:8:ARG:HD2	1:A:9:ARG:HB3	35	0.15
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB2	35	0.15
(1,64)	1:A:8:ARG:HD3	1:A:9:ARG:HB3	35	0.15
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	28	0.15
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	28	0.15
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	31	0.15
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	31	0.15
(1,54)	1:A:1:GLU:HA	1:A:2:LEU:HG	13	0.15
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	5	0.14
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	5	0.14
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	12	0.14
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	12	0.14
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	1	0.13
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	1	0.13
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	16	0.13
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	16	0.13
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	21	0.13
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	21	0.13
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	22	0.13
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	22	0.13
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	35	0.13
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	35	0.13
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	6	0.12
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	6	0.12
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	32	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	32	0.12
(1,54)	1:A:1:GLU:HA	1:A:2:LEU:HG	5	0.12
(1,54)	1:A:1:GLU:HA	1:A:2:LEU:HG	17	0.12
(1,54)	1:A:1:GLU:HA	1:A:2:LEU:HG	40	0.12
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD2	9	0.12
(1,25)	1:A:6:LYS:HA	1:A:6:LYS:HD3	9	0.12
(1,67)	1:A:9:ARG:HD2	1:A:8:ARG:HG2	4	0.11
(1,67)	1:A:9:ARG:HD2	1:A:8:ARG:HG3	4	0.11
(1,67)	1:A:9:ARG:HD3	1:A:8:ARG:HG2	4	0.11
(1,67)	1:A:9:ARG:HD3	1:A:8:ARG:HG3	4	0.11
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	2	0.11
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	2	0.11
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	3	0.11
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	3	0.11
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	8	0.11
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	8	0.11
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	15	0.11
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	15	0.11
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	20	0.11
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	20	0.11
(1,58)	1:A:7:PRO:HD2	1:A:9:ARG:HA	25	0.11
(1,58)	1:A:7:PRO:HD3	1:A:9:ARG:HA	25	0.11

## 10 Dihedral-angle violation analysis (i)

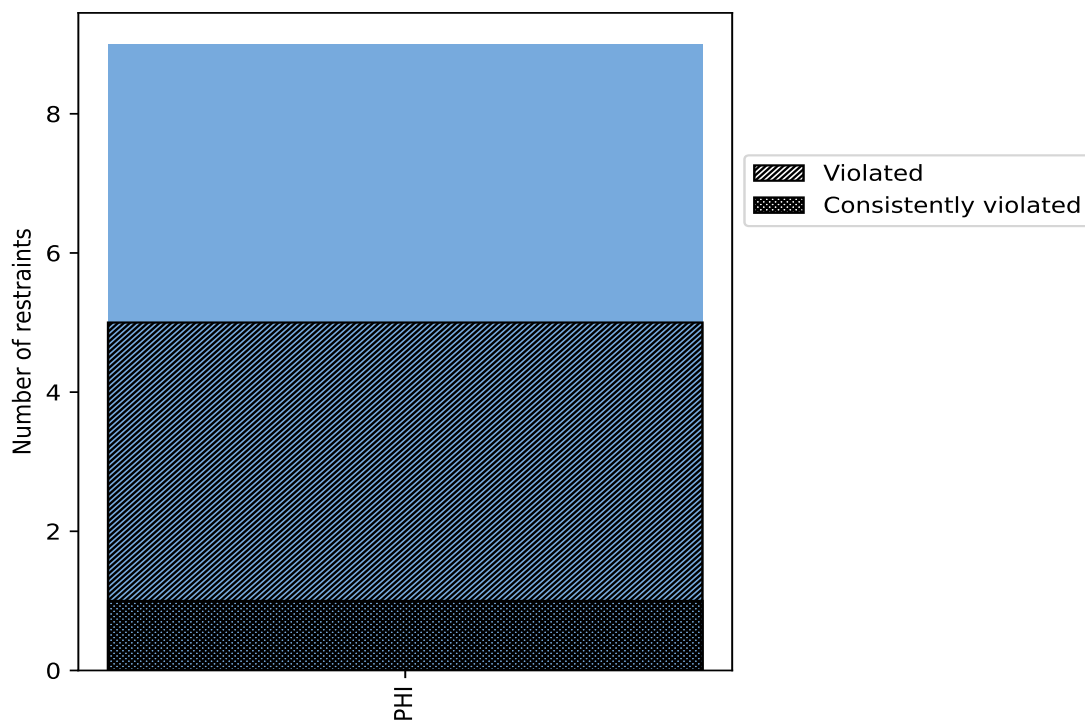
### 10.1 Summary of dihedral-angle violations (i)

The following table provides the summary of dihedral-angle violations in different dihedral-angle types. Violations less than 1° are not included in the calculation.

Angle 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>
PHI	9	100.0	5	55.6	55.6	1	11.1	11.1
Total	9	100.0	5	55.6	55.6	1	11.1	11.1

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

#### 10.1.1 Bar chart : Distribution of dihedral-angles and violations (i)



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

## 10.2 Dihedral-angle violation statistics for each model [\(i\)](#)

The following table provides the dihedral-angle violation statistics for each model in the ensemble. Violations less than 1° are not included in the statistics.

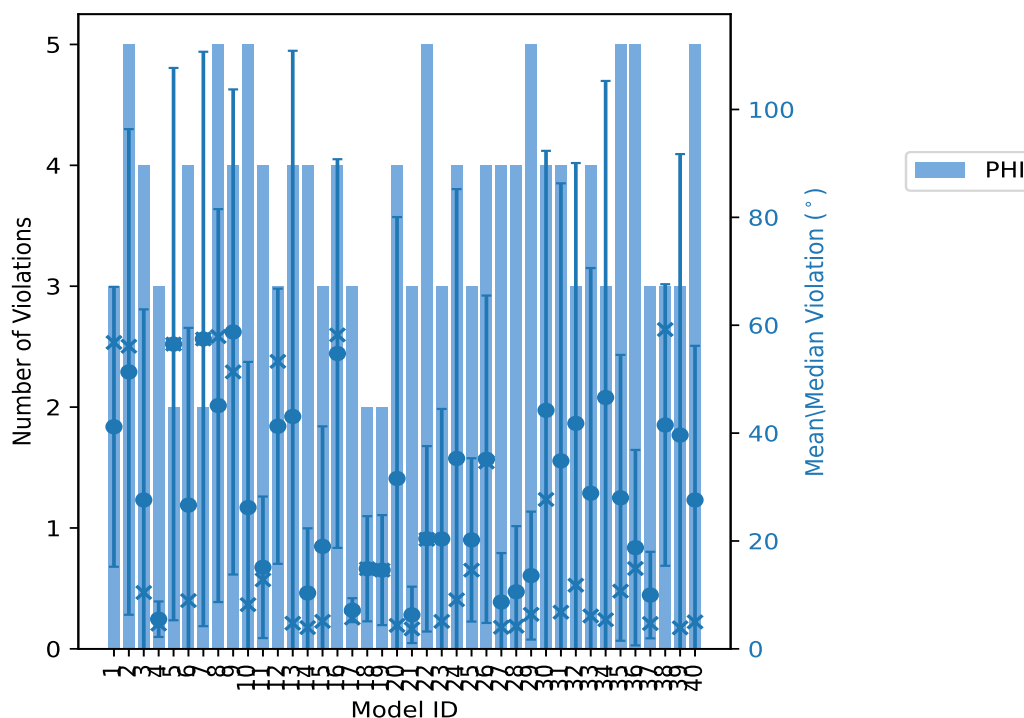
Model ID	Number of violations		Mean (°)	Max (°)	SD (°)	Median (°)
	PHI	Total				
1	3	3	41.17	62.1	25.95	56.8
2	5	5	51.34	123.6	45.02	56.1
3	4	4	27.6	88.1	35.34	10.45
4	3	3	5.5	9.9	3.29	4.6
5	2	2	56.5	107.7	51.2	56.5
6	4	4	26.65	83.5	32.86	8.95
7	2	2	57.45	110.7	53.25	57.45
8	5	5	45.1	94.9	36.43	57.9
9	4	4	58.75	127.2	44.97	51.35
10	5	5	26.2	68.8	27.0	8.2
11	4	4	15.12	33.5	13.12	12.8
12	3	3	41.27	64.7	25.51	53.3
13	4	4	43.08	160.5	67.8	4.75
14	4	4	10.35	31.1	12.0	3.95
15	3	3	19.0	50.4	22.25	5.1
16	4	4	54.75	97.6	36.03	58.2
17	3	3	7.17	10.3	2.22	5.7
18	2	2	14.85	24.6	9.75	14.85
19	2	2	14.6	24.8	10.2	14.6
20	4	4	31.58	115.5	48.47	4.35
21	3	3	6.3	13.6	5.23	3.7
22	5	5	20.4	49.0	17.19	20.3
23	3	3	20.37	54.4	24.11	5.1
24	4	4	35.3	121.5	49.94	9.1
25	3	3	20.2	40.9	15.14	14.6
26	4	4	35.15	67.6	30.35	34.65
27	4	4	8.68	24.3	9.08	4.0
28	4	4	10.58	31.6	12.17	4.2
29	5	5	13.58	32.8	11.86	6.4
30	4	4	44.22	120.4	48.11	27.7
31	4	4	34.85	123.9	51.46	6.8
32	3	3	41.8	109.9	48.27	11.8
33	4	4	28.85	101.1	41.75	6.1
34	3	3	46.6	129.6	58.69	5.4
35	5	5	28.0	63.2	26.5	10.7
36	5	5	18.76	52.3	18.12	14.9
37	3	3	9.97	21.3	8.02	4.7
38	3	3	41.5	60.7	26.1	59.2

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Model ID	Number of violations		Mean (°)	Max (°)	SD (°)	Median (°)
	PHI	Total				
39	3	3	39.63	113.3	52.1	3.9
40	5	5	27.6	63.0	28.58	5.0

### 10.2.1 Bar graph : Dihedral violation statistics for each model [\(i\)](#)



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

### 10.3 Dihedral-angle violation statistics for the ensemble [\(i\)](#)

Violation analysis may find that some restraints are violated in very 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 ensemble.

Number of violated restraints		Fraction of the ensemble	
PHI	Total	Count <sup>1</sup>	%
0	0	1	2.5
0	0	2	5.0
0	0	3	7.5
0	0	4	10.0

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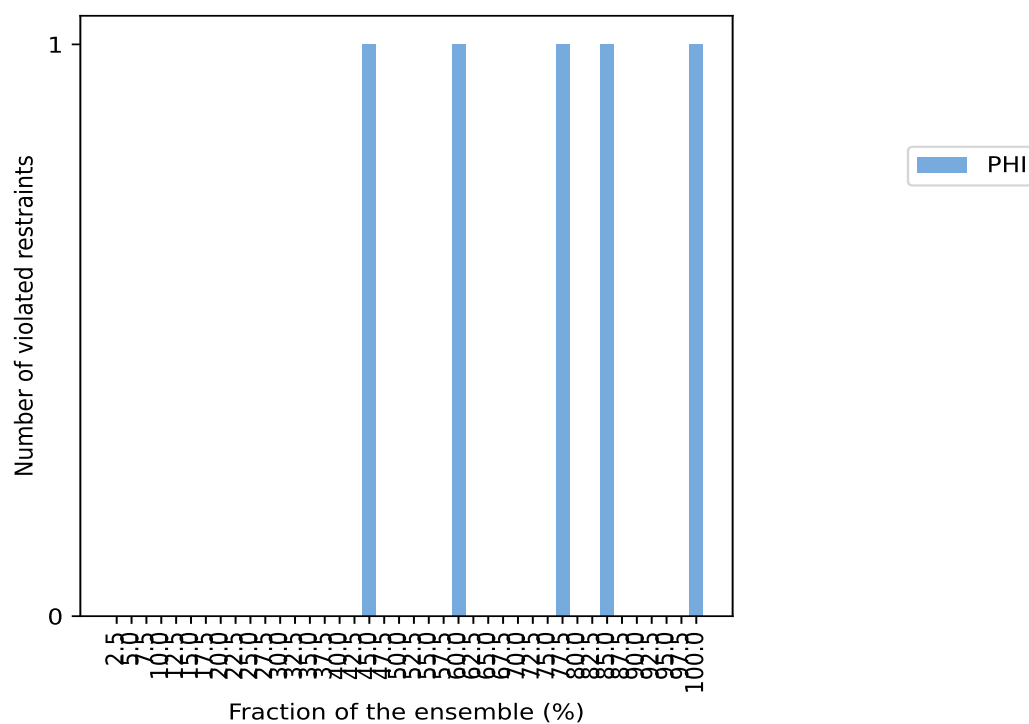


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PHI	Number of violated restraints		Fraction of the ensemble	
	Total	Count <sup>1</sup>	Count <sup>1</sup>	%
0	0	5	5	12.5
0	0	6	6	15.0
0	0	7	7	17.5
0	0	8	8	20.0
0	0	9	9	22.5
0	0	10	10	25.0
0	0	11	11	27.5
0	0	12	12	30.0
0	0	13	13	32.5
0	0	14	14	35.0
0	0	15	15	37.5
0	0	16	16	40.0
0	0	17	17	42.5
1	1	18	18	45.0
0	0	19	19	47.5
0	0	20	20	50.0
0	0	21	21	52.5
0	0	22	22	55.0
0	0	23	23	57.5
1	1	24	24	60.0
0	0	25	25	62.5
0	0	26	26	65.0
0	0	27	27	67.5
0	0	28	28	70.0
0	0	29	29	72.5
0	0	30	30	75.0
1	1	31	31	77.5
0	0	32	32	80.0
0	0	33	33	82.5
1	1	34	34	85.0
0	0	35	35	87.5
0	0	36	36	90.0
0	0	37	37	92.5
0	0	38	38	95.0
0	0	39	39	97.5
1	1	40	40	100.0

<sup>1</sup> Number of models with violations

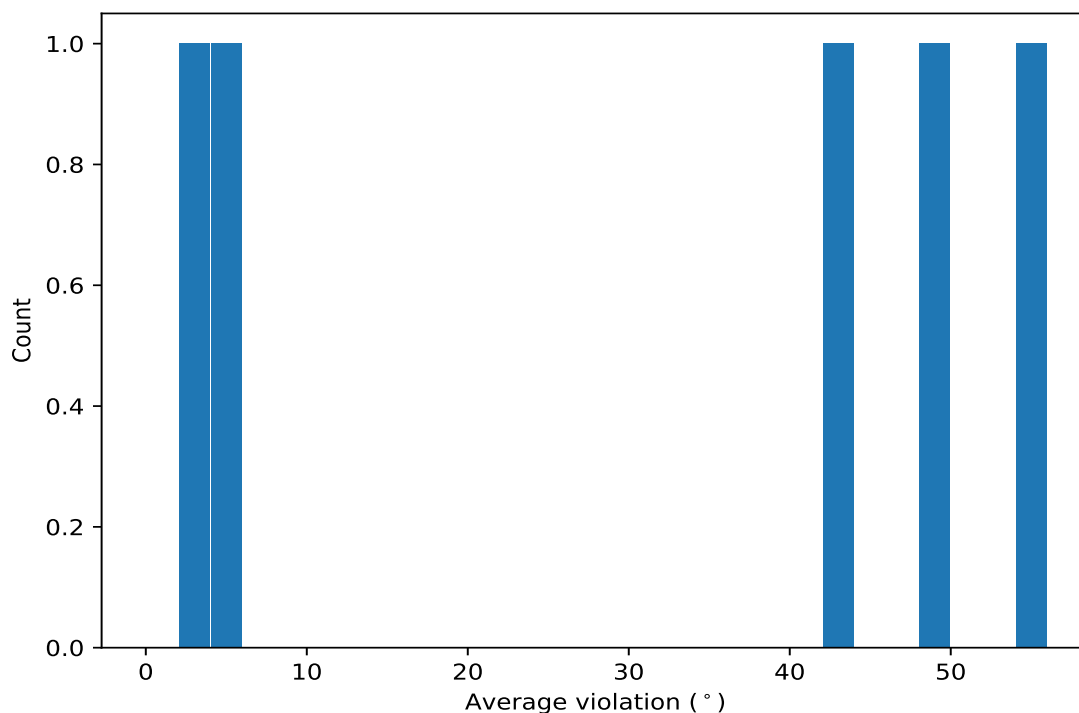
### 10.3.1 Bar graph : Dihedral-angle Violation statistics for the ensemble [i](#)



## 10.4 Most violated dihedral-angle restraints in the ensemble [i](#)

### 10.4.1 Histogram : Distribution of mean dihedral-angle 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



#### 10.4.2 Table: Most violated dihedral-angle 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.

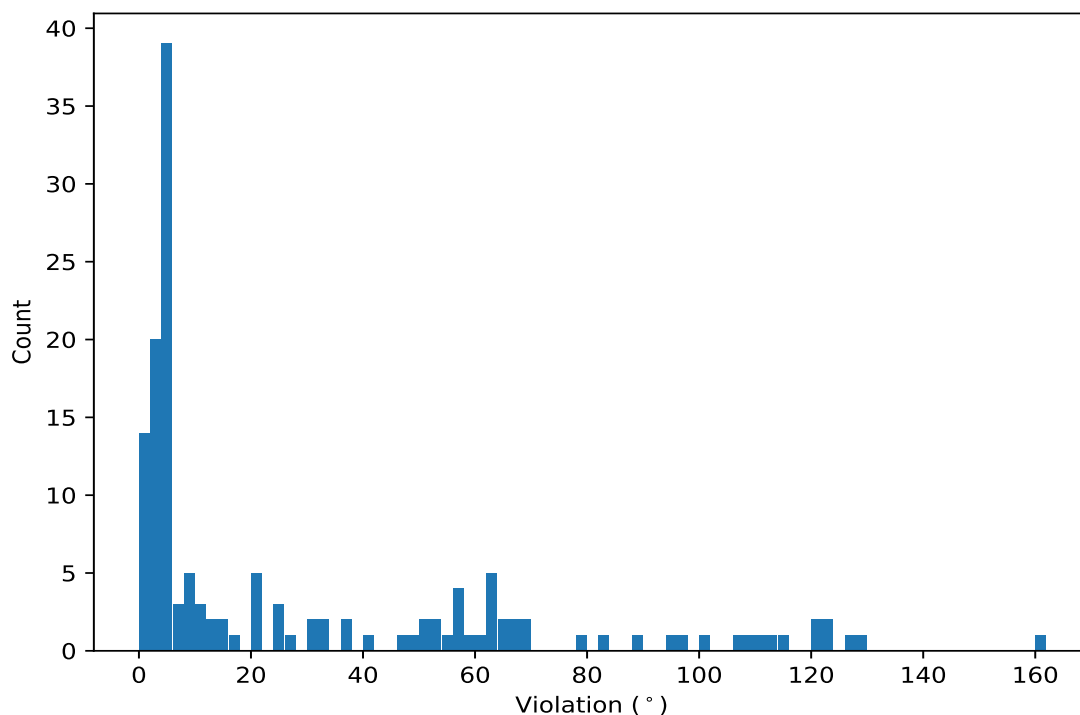
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models <sup>1</sup>	Mean	SD <sup>2</sup>	Median
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	40	4.57	0.87	4.95
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	34	55.41	37.44	55.6
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	31	42.84	47.21	16.0
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	24	2.22	0.98	1.9
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	18	48.73	24.08	54.7

<sup>1</sup> Number of violated models, <sup>2</sup>Standard deviation, All angle values are in degree (°)

## 10.5 All violated dihedral-angle restraints [i](#)

### 10.5.1 Histogram : Distribution of violations [i](#)

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



### 10.5.2 Table: All violated dihedral-angle restraints [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.

Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	13	160.5
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	34	129.6
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	9	127.2
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	31	123.9
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	2	123.6
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	24	121.5
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	30	120.4
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	20	115.5
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	39	113.3
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	7	110.7
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	32	109.9
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	5	107.7
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	33	101.1
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	16	97.6
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	8	94.9
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	3	88.1
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	6	83.5
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	16	78.9
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	2	69.7
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	10	68.8
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	26	67.6

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	8	66.5
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	9	65.8
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	12	64.7
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	26	63.3
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	35	63.2
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	40	63.0
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	40	62.2
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	1	62.1
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	38	60.7
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	38	59.2
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	8	57.9
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	35	57.3
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	1	56.8
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	2	56.1
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	23	54.4
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	12	53.3
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	36	52.3
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	30	50.6
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	15	50.4
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	22	49.0
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	10	47.4
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	25	40.9
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	16	37.5
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	9	36.9
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	11	33.5
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	29	32.8
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	28	31.6
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	14	31.1
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	22	27.0
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	19	24.8
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	18	24.6
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	27	24.3
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	29	21.9
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	11	21.5
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	37	21.3
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	36	20.5
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	22	20.3
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	3	16.0
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	36	14.9
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	25	14.6
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	21	13.6
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	24	12.9
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	32	11.8
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	35	10.7
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	17	10.3
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	4	9.9
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	6	9.2
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	6	8.7
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	31	8.4
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	10	8.2
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	33	7.3

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	29	6.4
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	26	6.0
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	12	5.8
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	17	5.7
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	13	5.5
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	17	5.5
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	2	5.4
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	34	5.4
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	5	5.3
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	10	5.3
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	24	5.3
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	27	5.2
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	6	5.2
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	31	5.2
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	9	5.1
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	15	5.1
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	18	5.1
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	23	5.1
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	25	5.1
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	29	5.1
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	8	5.0
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	16	5.0
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	20	5.0
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	35	5.0
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	40	5.0
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	3	4.9
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	33	4.9
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	34	4.8
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	30	4.8
(1,8)	1:A:11:TYR:C	1:A:12:ILE:N	1:A:12:ILE:CA	1:A:12:ILE:C	28	4.7
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	37	4.7
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	1	4.6
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	4	4.6
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	36	4.6
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	38	4.6
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	40	4.5
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	19	4.4
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	7	4.2
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	11	4.1
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	14	4.1
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	13	4.0
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	37	3.9
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	39	3.9
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	14	3.8
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	35	3.8
(1,7)	1:A:10:PRO:C	1:A:11:TYR:N	1:A:11:TYR:CA	1:A:11:TYR:C	20	3.7
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	28	3.7
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	21	3.7
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	26	3.7
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	32	3.7
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	40	3.3

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*Continued from previous page...*

Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	22	3.0
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	27	2.8
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	22	2.7
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	14	2.4
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	27	2.4
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	13	2.3
(1,5)	1:A:7:PRO:C	1:A:8:ARG:N	1:A:8:ARG:CA	1:A:8:ARG:C	28	2.3
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	20	2.1
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	33	2.1
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	4	2.0
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	2	1.9
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	31	1.9
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	29	1.7
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	39	1.7
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	21	1.6
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	23	1.6
(1,9)	1:A:12:ILE:C	1:A:13:LEU:N	1:A:13:LEU:CA	1:A:13:LEU:C	15	1.5
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	24	1.5
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	36	1.5
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	3	1.4
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	11	1.4
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	10	1.3
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	8	1.2
(1,6)	1:A:8:ARG:C	1:A:9:ARG:N	1:A:9:ARG:CA	1:A:9:ARG:C	30	1.1