



Full wwPDB NMR Structure Validation Report ⓘ

Apr 29, 2025 – 06:31 PM EDT

PDB ID : 9BF4 / pdb_00009bf4
BMRB ID : 31168
Title : NMR solution structure of Paws Derived Peptide-25 (PDP-25)
Authors : Hajiaghaalipour, F.; Wong, W.; Payne, C.D.; Fisher, M.F.; Clark, R.J.; Mylne, J.S.; Rosengren, K.J.
Deposited on : 2024-04-16

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at 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>.

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

MolProbity : 4-5-2 with Phenix2.0rc1
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) : 2.43.1

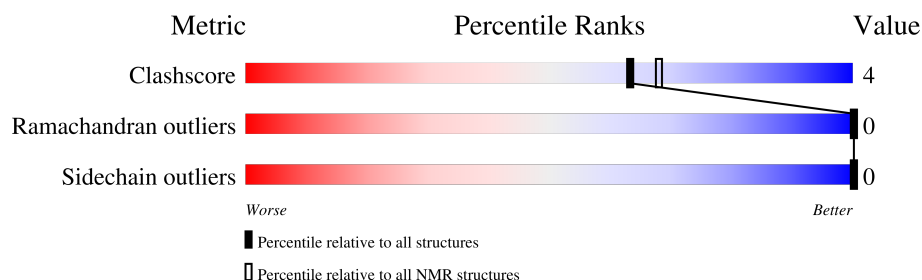
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 73%.

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 ≥ 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	31	<div> <div style="width: 94%;"></div> <div style="width: 6%;"></div> </div> <div>94%6%</div>

2 Ensemble composition and analysis

This entry contains 20 models. Model 2 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:1-A:31 (31)	0.75	2

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 3 clusters. No single-model clusters were found.

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

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 438 atoms, of which 204 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Paws Derived Peptide-25.

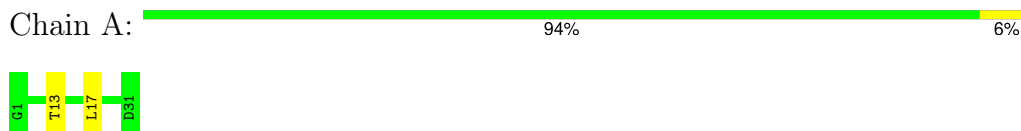
Mol	Chain	Residues	Atoms						Trace
1	A	31	Total	C	H	N	O	S	0
			438	151	204	35	43	5	

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: Paws Derived Peptide-25

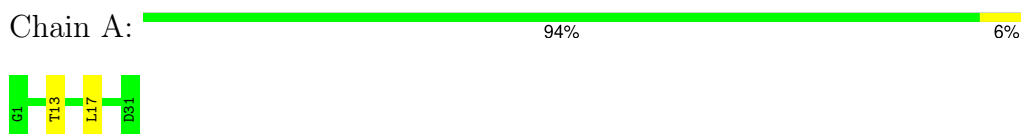


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: Paws Derived Peptide-25



4.2.2 Score per residue for model 2 (medoid)

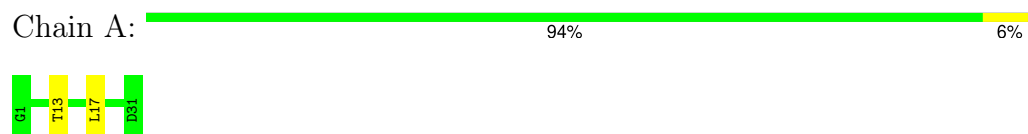
- Molecule 1: Paws Derived Peptide-25



There are no outlier residues in this chain.

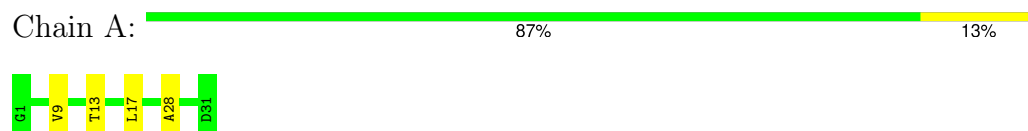
4.2.3 Score per residue for model 3

- Molecule 1: Paws Derived Peptide-25



4.2.4 Score per residue for model 4

- Molecule 1: Paws Derived Peptide-25



4.2.5 Score per residue for model 5

- Molecule 1: Paws Derived Peptide-25



There are no outlier residues in this chain.

4.2.6 Score per residue for model 6

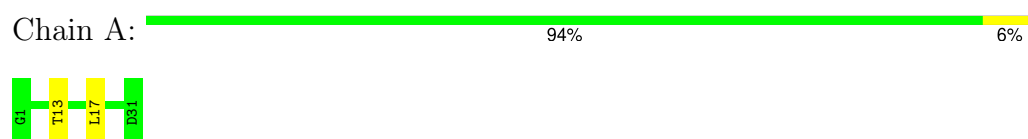
- Molecule 1: Paws Derived Peptide-25



There are no outlier residues in this chain.

4.2.7 Score per residue for model 7

- Molecule 1: Paws Derived Peptide-25



4.2.8 Score per residue for model 8

- Molecule 1: Paws Derived Peptide-25





4.2.9 Score per residue for model 9

- Molecule 1: Paws Derived Peptide-25

Chain A:
81% 19%



4.2.10 Score per residue for model 10

- Molecule 1: Paws Derived Peptide-25

Chain A:
94% 6%



4.2.11 Score per residue for model 11

- Molecule 1: Paws Derived Peptide-25

Chain A:
87% 13%



4.2.12 Score per residue for model 12

- Molecule 1: Paws Derived Peptide-25

Chain A:
94% 6%



4.2.13 Score per residue for model 13

- Molecule 1: Paws Derived Peptide-25

Chain A:
87% 13%



4.2.14 Score per residue for model 14

- Molecule 1: Paws Derived Peptide-25

Chain A: 90% 10%



4.2.15 Score per residue for model 15

- Molecule 1: Paws Derived Peptide-25

Chain A: 77% 23%



4.2.16 Score per residue for model 16

- Molecule 1: Paws Derived Peptide-25

Chain A: 87% 13%



4.2.17 Score per residue for model 17

- Molecule 1: Paws Derived Peptide-25

Chain A: 77% 19% 4%



4.2.18 Score per residue for model 18

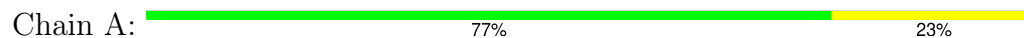
- Molecule 1: Paws Derived Peptide-25

Chain A: 87% 13%



4.2.19 Score per residue for model 19

- Molecule 1: Paws Derived Peptide-25



4.2.20 Score per residue for model 20

- Molecule 1: Paws Derived Peptide-25



5 Refinement protocol and experimental data overview

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

Of the 50 calculated structures, 20 were deposited, based on the following criterion: *structures with acceptable covalent geometry*.

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

Software name	Classification	Version
CYANA	structure calculation	
CNS	refinement	1.21

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	275
Number of shifts mapped to atoms	275
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	73%

6 Model quality [i](#)

6.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	1.00±0.05	0±0/243 (0.0± 0.0%)	0.93±0.03	0±0/333 (0.0± 0.1%)
All	All	1.00	0/4860 (0.0%)	0.93	2/6660 (0.0%)

There are no bond-length outliers.

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	18	PRO	CA-C-N	5.27	126.42	119.84	17	1
1	A	18	PRO	C-N-CA	5.27	126.42	119.84	17	1

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	234	204	204	2±1
All	All	4680	4080	4080	35

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 4.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:1:GLY:HA2	1:A:28:ALA:O	0.56	2.01	17	4
1:A:13:THR:O	1:A:17:LEU:HG	0.55	2.01	8	15
1:A:14:CYS:O	1:A:22:GLN:HG2	0.54	2.01	19	5
1:A:10:PRO:HB2	1:A:13:THR:OG1	0.51	2.06	14	4
1:A:18:PRO:O	1:A:22:GLN:HG3	0.47	2.09	17	2
1:A:13:THR:O	1:A:17:LEU:HD13	0.47	2.09	11	2
1:A:9:VAL:HG11	1:A:28:ALA:HB1	0.41	1.93	4	2
1:A:14:CYS:N	1:A:25:CYS:SG	0.40	2.95	16	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	29/31 (94%)	28±1 (97±2%)	1±1 (3±2%)	0±0 (0±0%)	100	100
All	All	580/620 (94%)	561 (97%)	19 (3%)	0 (0%)	100	100

There are no Ramachandran outliers.

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	25/25 (100%)	25±0 (100±0%)	0±0 (0±0%)	100	100
All	All	500/500 (100%)	500 (100%)	0 (0%)	100	100

There are no protein residues with a non-rotameric sidechain to report.

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 73% for the well-defined parts and 73% for the entire structure.

7.1 Chemical shift list 1

File name: `working_cs.cif`

Chemical shift list name: *starch_output*

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	275
Number of shifts mapped to atoms	275
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

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	30	3.10 ± 0.13	Should be applied
$^{13}\text{C}_\beta$	28	3.16 ± 0.26	Should be applied
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	27	1.99 ± 0.28	Should be applied

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 73%, i.e. 274 atoms were assigned a chemical shift out of a possible 374. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	116/152 (76%)	59/62 (95%)	30/62 (48%)	27/28 (96%)
Sidechain	158/169 (93%)	108/112 (96%)	50/55 (91%)	0/2 (0%)

Continued on next page...

Continued from previous page...

	Total	¹ H	¹³ C	¹⁵ N
Aromatic	0/53 (0%)	0/26 (0%)	0/25 (0%)	0/2 (0%)
Overall	274/374 (73%)	167/200 (84%)	80/142 (56%)	27/32 (84%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 73%, i.e. 274 atoms were assigned a chemical shift out of a possible 374. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	116/152 (76%)	59/62 (95%)	30/62 (48%)	27/28 (96%)
Sidechain	158/169 (93%)	108/112 (96%)	50/55 (91%)	0/2 (0%)
Aromatic	0/53 (0%)	0/26 (0%)	0/25 (0%)	0/2 (0%)
Overall	274/374 (73%)	167/200 (84%)	80/142 (56%)	27/32 (84%)

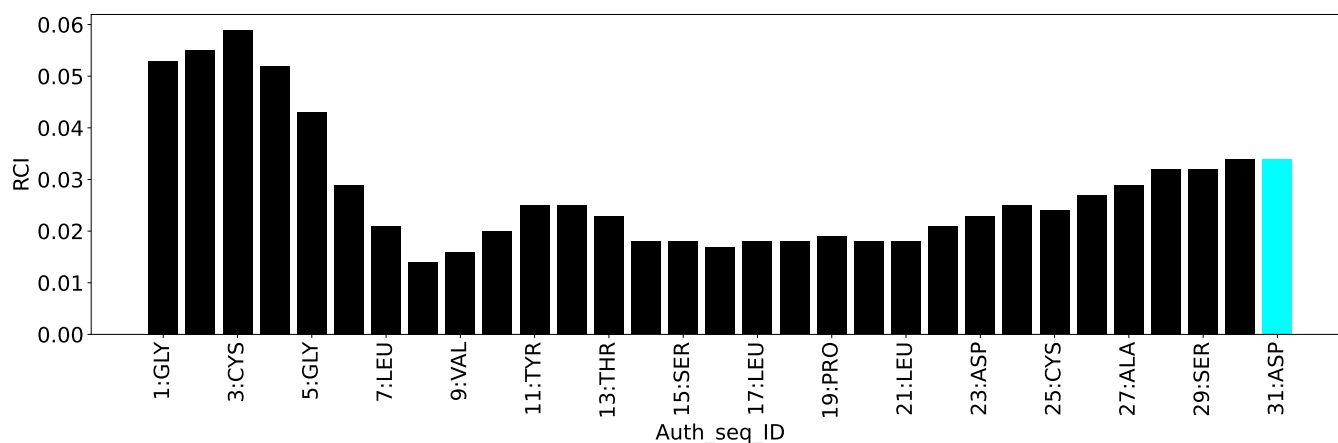
7.1.4 Statistically unusual chemical shifts [i](#)

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	348
Intra-residue ($ i-j =0$)	104
Sequential ($ i-j =1$)	121
Medium range ($ i-j >1$ and $ i-j <5$)	63
Long range ($ i-j \geq 5$)	28
Inter-chain	0
Hydrogen bond restraints	32
Disulfide bond restraints	0
Total dihedral-angle restraints	63
Number of unmapped restraints	0
Number of restraints per residue	13.3
Number of long range restraints per residue ¹	1.0

¹Long 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)	3.0	0.2
0.2-0.5 (Medium)	0.6	0.42
>0.5 (Large)	0.8	1.58

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)	0.4	1.95
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

9 Distance violation analysis ⓘ

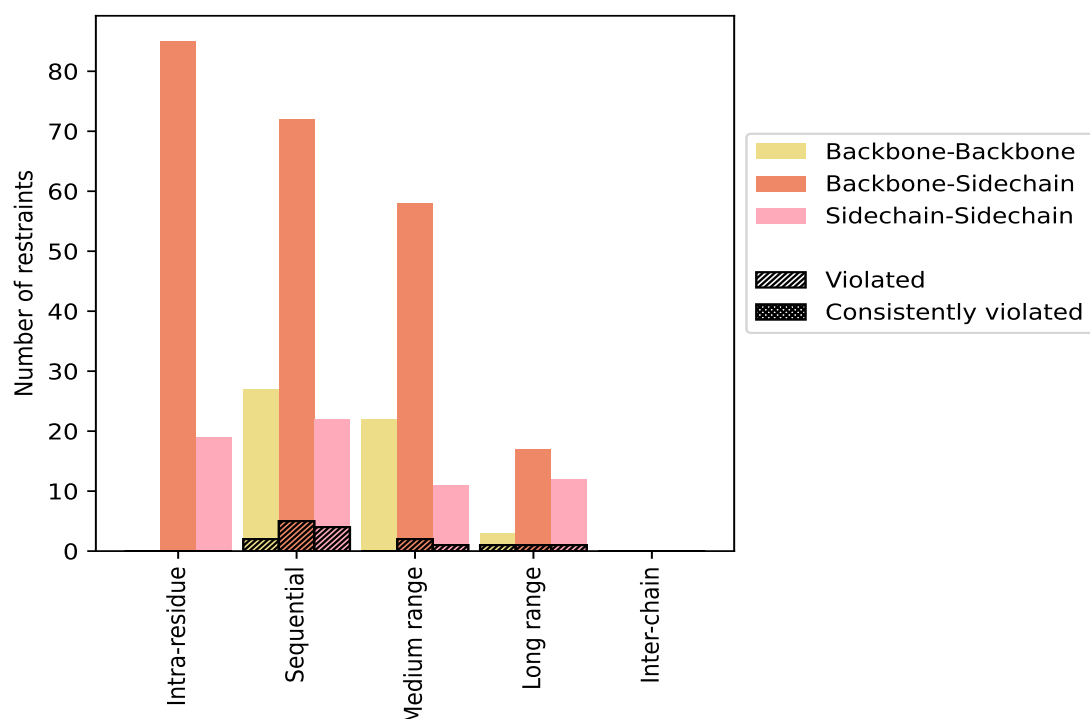
9.1 Summary of distance violations ⓘ

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	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	104	29.9	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	85	24.4	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	19	5.5	0	0.0	0.0	0	0.0	0.0
Sequential ($i-j =1$)	121	34.8	11	9.1	3.2	0	0.0	0.0
Backbone-Backbone	27	7.8	2	7.4	0.6	0	0.0	0.0
Backbone-Sidechain	72	20.7	5	6.9	1.4	0	0.0	0.0
Sidechain-Sidechain	22	6.3	4	18.2	1.1	0	0.0	0.0
Medium range ($i-j >1$ & $i-j <5$)	63	18.1	2	3.2	0.6	0	0.0	0.0
Backbone-Backbone	22	6.3	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	30	8.6	1	3.3	0.3	0	0.0	0.0
Sidechain-Sidechain	11	3.2	1	9.1	0.3	0	0.0	0.0
Long range ($i-j \geq 5$)	28	8.0	2	7.1	0.6	0	0.0	0.0
Backbone-Backbone	3	0.9	1	33.3	0.3	0	0.0	0.0
Backbone-Sidechain	13	3.7	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	12	3.4	1	8.3	0.3	0	0.0	0.0
Inter-chain	0	0.0	0	0.0	0.0	0	0.0	0.0
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
Hydrogen bond	32	9.2	2	6.2	0.6	0	0.0	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	348	100.0	17	4.9	4.9	0	0.0	0.0
Backbone-Backbone	52	14.9	3	5.8	0.9	0	0.0	0.0
Backbone-Sidechain	232	66.7	8	3.4	2.3	0	0.0	0.0
Sidechain-Sidechain	64	18.4	6	9.4	1.7	0	0.0	0.0

¹ percentage calculated with respect to the total number of distance restraints, ² percentage calculated with respect to the number of restraints in a particular restraint category, ³ violated in at least one model, ⁴ 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 ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	0	2	0	1	0	3	0.45	0.68	0.25	0.58
2	0	5	0	1	0	6	0.15	0.2	0.03	0.15
3	0	4	0	1	0	5	0.33	1.01	0.34	0.17
4	0	4	0	2	0	6	0.17	0.33	0.08	0.14
5	0	3	0	1	0	4	0.57	1.11	0.42	0.53
6	0	4	1	1	0	6	0.15	0.29	0.06	0.12
7	0	2	0	2	0	4	0.26	0.51	0.15	0.22
8	0	5	0	2	0	7	0.43	1.44	0.49	0.13
9	0	5	0	1	0	6	0.37	1.08	0.38	0.13
10	0	4	0	1	0	5	0.48	1.08	0.42	0.19

Continued on next page...

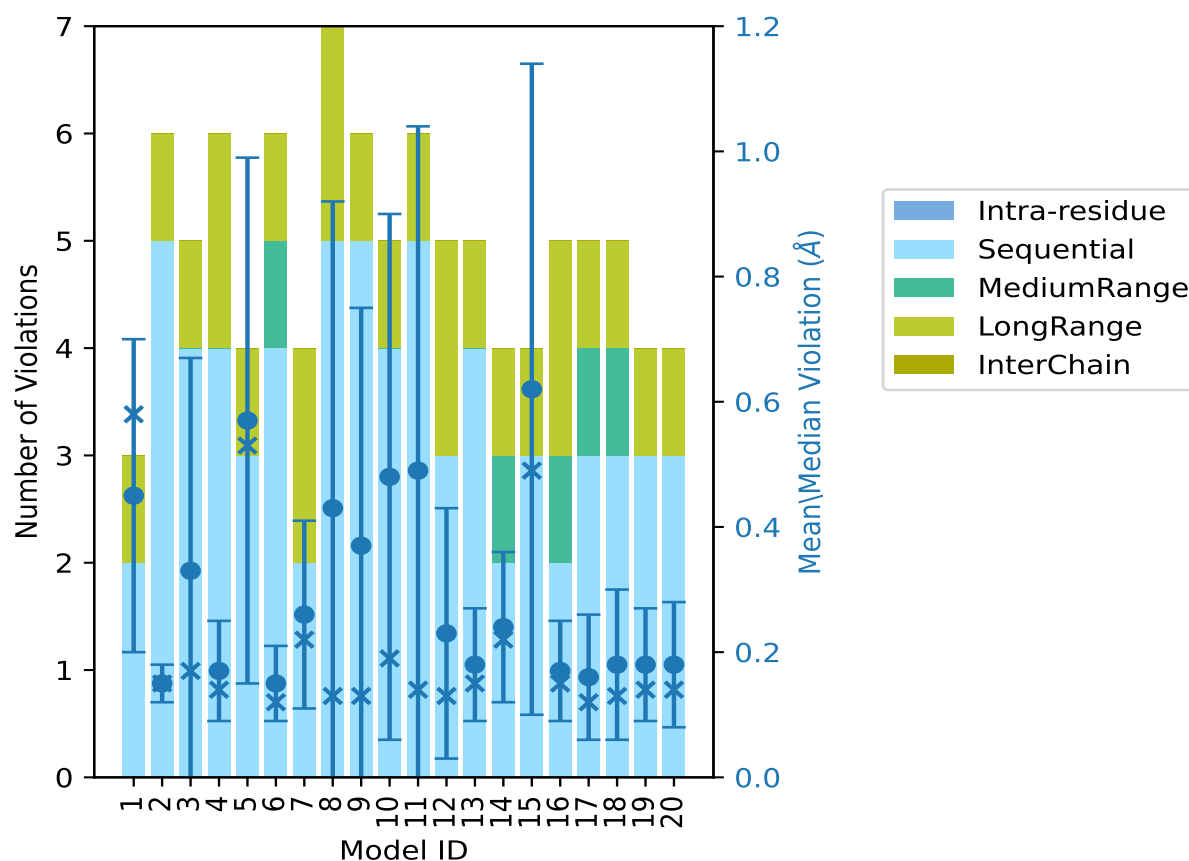
Continued from previous page...

Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
11	0	5	0	1	0	6	0.49	1.58	0.55	0.14
12	0	3	0	2	0	5	0.23	0.62	0.2	0.13
13	0	4	0	1	0	5	0.18	0.36	0.09	0.15
14	0	2	1	1	0	4	0.24	0.41	0.12	0.22
15	0	3	0	1	0	4	0.62	1.39	0.52	0.49
16	0	2	1	2	0	5	0.17	0.32	0.08	0.15
17	0	3	1	1	0	5	0.16	0.36	0.1	0.12
18	0	3	1	1	0	5	0.18	0.42	0.12	0.13
19	0	3	0	1	0	4	0.18	0.34	0.09	0.14
20	0	3	0	1	0	4	0.18	0.36	0.1	0.14

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints,

⁵Inter-chain restraints, ⁶Standard 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

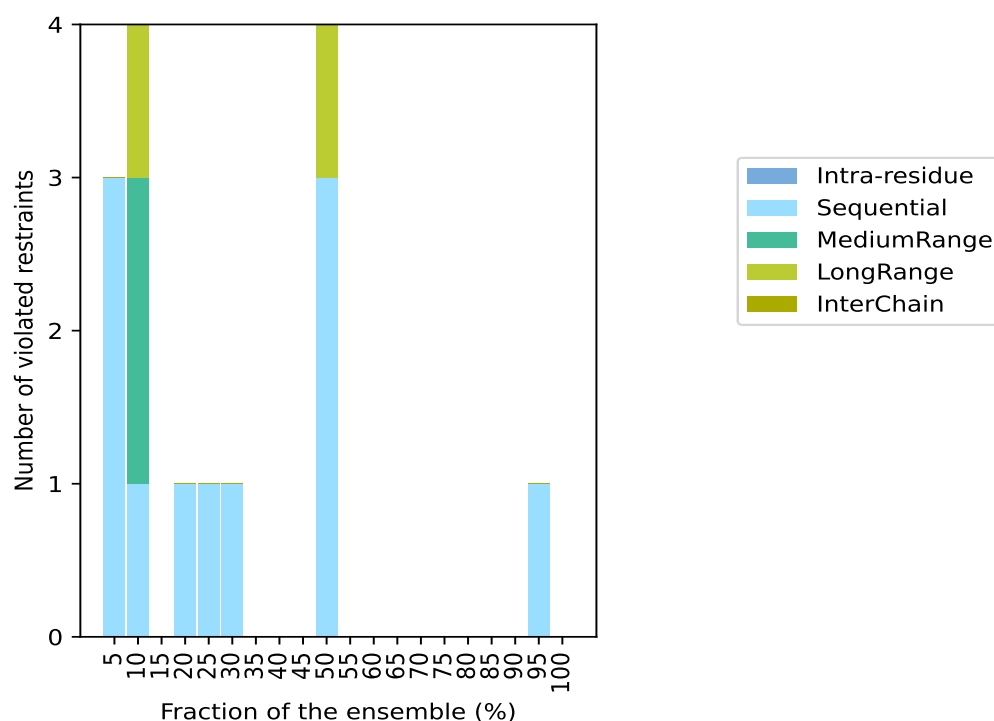
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 301(IR:104, SQ:110, MR:61, LR:26, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
0	3	0	0	0	3	1	5.0
0	1	2	1	0	4	2	10.0
0	0	0	0	0	0	3	15.0
0	1	0	0	0	1	4	20.0
0	1	0	0	0	1	5	25.0
0	1	0	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	3	0	1	0	4	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	1	0	0	0	1	19	95.0
0	0	0	0	0	0	20	100.0

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints,

⁵Inter-chain restraints, ⁶ Number of models with violations

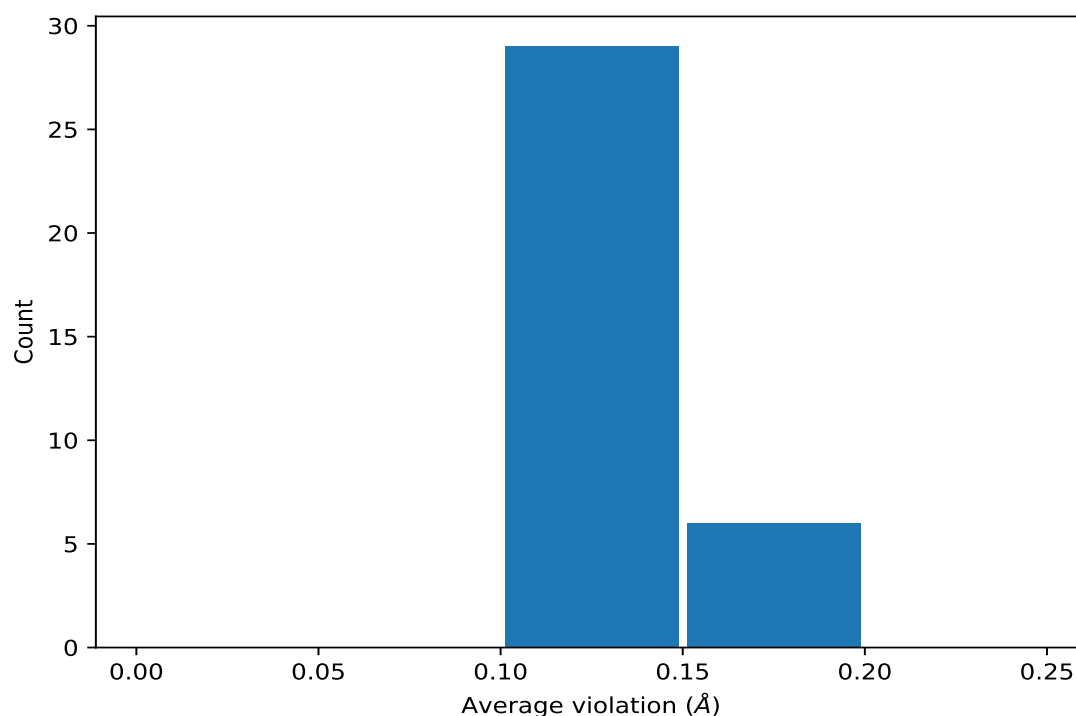
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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 ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	19	0.16	0.02	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	19	0.16	0.02	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	19	0.16	0.02	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	19	0.16	0.02	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	19	0.16	0.02	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	19	0.16	0.02	0.16
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	10	0.12	0.01	0.12
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	10	0.12	0.02	0.11
(1,196)	1:9:A:VAL:HB	1:10:A:PRO:HD3	6	0.11	0.01	0.11
(1,55)	1:3:A:CYS:HB2	1:4:A:TRP:H	5	0.11	0.01	0.11
(1,227)	1:4:A:TRP:HD1	1:5:A:GLY:H	4	0.12	0.0	0.12
(1,115)	1:9:A:VAL:HG11	1:13:A:THR:H	2	0.12	0.02	0.12
(1,115)	1:9:A:VAL:HG12	1:13:A:THR:H	2	0.12	0.02	0.12
(1,115)	1:9:A:VAL:HG13	1:13:A:THR:H	2	0.12	0.02	0.12
(1,115)	1:9:A:VAL:HG21	1:13:A:THR:H	2	0.12	0.02	0.12
(1,115)	1:9:A:VAL:HG22	1:13:A:THR:H	2	0.12	0.02	0.12

Continued on next page...

Continued from previous page...

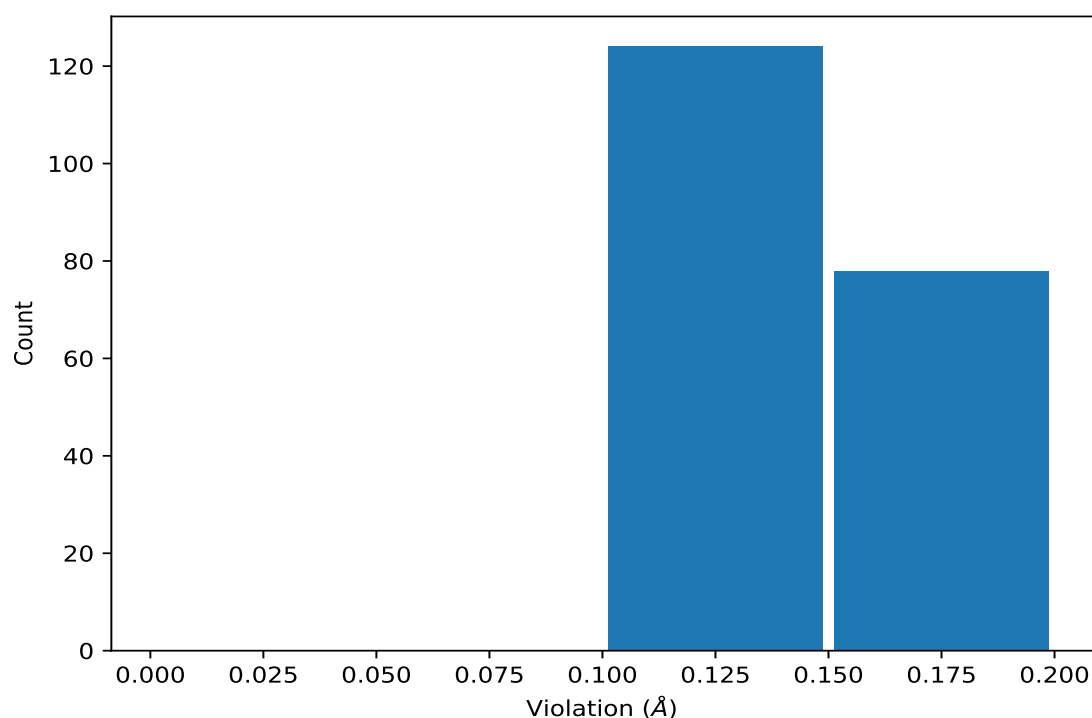
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,115)	1:9:A:VAL:HG23	1:13:A:THR:H	2	0.12	0.02	0.12
(1,221)	1:10:A:PRO:HG2	1:11:A:TYR:HD1	2	0.12	0.0	0.12
(1,221)	1:10:A:PRO:HG2	1:11:A:TYR:HD2	2	0.12	0.0	0.12
(1,221)	1:10:A:PRO:HG3	1:11:A:TYR:HD1	2	0.12	0.0	0.12
(1,221)	1:10:A:PRO:HG3	1:11:A:TYR:HD2	2	0.12	0.0	0.12
(1,274)	1:17:A:LEU:HD11	1:22:A:GLN:HE21	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD11	1:22:A:GLN:HE22	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD12	1:22:A:GLN:HE21	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD12	1:22:A:GLN:HE22	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD13	1:22:A:GLN:HE21	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD13	1:22:A:GLN:HE22	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD21	1:22:A:GLN:HE21	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD21	1:22:A:GLN:HE22	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD22	1:22:A:GLN:HE21	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD22	1:22:A:GLN:HE22	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD23	1:22:A:GLN:HE21	2	0.11	0.0	0.11
(1,274)	1:17:A:LEU:HD23	1:22:A:GLN:HE22	2	0.11	0.0	0.11
(1,283)	1:18:A:PRO:HG2	1:22:A:GLN:HE21	2	0.11	0.0	0.11
(1,283)	1:18:A:PRO:HG2	1:22:A:GLN:HE22	2	0.11	0.0	0.11

¹Number of violated models, ²Standard 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)
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	2	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	2	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	2	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	2	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	2	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	2	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	5	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	5	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	5	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	5	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	5	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	5	0.2
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	10	0.19
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	10	0.19
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	10	0.19
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	10	0.19

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	10	0.19
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	10	0.19
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	12	0.18
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	12	0.18
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	12	0.18
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	12	0.18
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	12	0.18
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	12	0.18
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	3	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	3	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	3	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	3	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	3	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	3	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	8	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	8	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	8	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	8	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	8	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	8	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	15	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	15	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	15	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	15	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	15	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	15	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	16	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	16	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	16	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	16	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	16	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	16	0.17
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	6	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	6	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	6	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	6	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	6	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	6	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	7	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	7	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	7	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	7	0.16

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	7	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	7	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	13	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	13	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	13	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	13	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	13	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	13	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	19	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	19	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	19	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	19	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	19	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	19	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	20	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	20	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	20	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	20	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	20	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	20	0.16
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	4	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	4	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	4	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	4	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	4	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	4	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	9	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	9	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	9	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	9	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	9	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	9	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	11	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	11	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	11	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	11	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	11	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	11	0.15
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	2	0.15
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	13	0.15
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	18	0.14
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	18	0.14

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	18	0.14
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	18	0.14
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	18	0.14
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	18	0.14
(1,115)	1:9:A:VAL:HG11	1:13:A:THR:H	14	0.14
(1,115)	1:9:A:VAL:HG12	1:13:A:THR:H	14	0.14
(1,115)	1:9:A:VAL:HG13	1:13:A:THR:H	14	0.14
(1,115)	1:9:A:VAL:HG21	1:13:A:THR:H	14	0.14
(1,115)	1:9:A:VAL:HG22	1:13:A:THR:H	14	0.14
(1,115)	1:9:A:VAL:HG23	1:13:A:THR:H	14	0.14
(1,80)	1:29:A:SER:HB3	1:30:A:PHE:H	3	0.14
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	14	0.13
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	14	0.13
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	14	0.13
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	14	0.13
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	14	0.13
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	14	0.13
(1,196)	1:9:A:VAL:HB	1:10:A:PRO:HD3	18	0.13
(1,180)	1:11:A:TYR:HD1	1:12:A:GLY:H	8	0.13
(1,180)	1:11:A:TYR:HD2	1:12:A:GLY:H	8	0.13
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	6	0.13
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	11	0.13
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	4	0.13
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	8	0.13
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	12	0.13
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	13	0.13
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	15	0.13
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD11	17	0.12
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD12	17	0.12
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD13	17	0.12
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD21	17	0.12
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD22	17	0.12
(2,1)	1:20:A:TRP:HE3	1:21:A:LEU:HD23	17	0.12
(1,227)	1:4:A:TRP:HD1	1:5:A:GLY:H	3	0.12
(1,227)	1:4:A:TRP:HD1	1:5:A:GLY:H	6	0.12
(1,227)	1:4:A:TRP:HD1	1:5:A:GLY:H	10	0.12
(1,221)	1:10:A:PRO:HG2	1:11:A:TYR:HD1	4	0.12
(1,221)	1:10:A:PRO:HG2	1:11:A:TYR:HD2	4	0.12
(1,221)	1:10:A:PRO:HG3	1:11:A:TYR:HD1	4	0.12
(1,221)	1:10:A:PRO:HG3	1:11:A:TYR:HD2	4	0.12
(1,196)	1:9:A:VAL:HB	1:10:A:PRO:HD3	2	0.12
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	5	0.12

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	11	0.12
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	17	0.12
(1,55)	1:3:A:CYS:HB2	1:4:A:TRP:H	19	0.12
(1,283)	1:18:A:PRO:HG2	1:22:A:GLN:HE21	18	0.11
(1,283)	1:18:A:PRO:HG2	1:22:A:GLN:HE22	18	0.11
(1,274)	1:17:A:LEU:HD11	1:22:A:GLN:HE21	8	0.11
(1,274)	1:17:A:LEU:HD11	1:22:A:GLN:HE22	8	0.11
(1,274)	1:17:A:LEU:HD12	1:22:A:GLN:HE21	8	0.11
(1,274)	1:17:A:LEU:HD12	1:22:A:GLN:HE22	8	0.11
(1,274)	1:17:A:LEU:HD13	1:22:A:GLN:HE21	8	0.11
(1,274)	1:17:A:LEU:HD13	1:22:A:GLN:HE22	8	0.11
(1,274)	1:17:A:LEU:HD21	1:22:A:GLN:HE21	8	0.11
(1,274)	1:17:A:LEU:HD21	1:22:A:GLN:HE22	8	0.11
(1,274)	1:17:A:LEU:HD22	1:22:A:GLN:HE21	8	0.11
(1,274)	1:17:A:LEU:HD22	1:22:A:GLN:HE22	8	0.11
(1,274)	1:17:A:LEU:HD23	1:22:A:GLN:HE21	8	0.11
(1,274)	1:17:A:LEU:HD23	1:22:A:GLN:HE22	8	0.11
(1,227)	1:4:A:TRP:HD1	1:5:A:GLY:H	12	0.11
(1,221)	1:10:A:PRO:HG2	1:11:A:TYR:HD1	11	0.11
(1,221)	1:10:A:PRO:HG2	1:11:A:TYR:HD2	11	0.11
(1,221)	1:10:A:PRO:HG3	1:11:A:TYR:HD1	11	0.11
(1,221)	1:10:A:PRO:HG3	1:11:A:TYR:HD2	11	0.11
(1,196)	1:9:A:VAL:HB	1:10:A:PRO:HD3	20	0.11
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	7	0.11
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	16	0.11
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	20	0.11
(1,115)	1:9:A:VAL:HG11	1:13:A:THR:H	16	0.11
(1,115)	1:9:A:VAL:HG12	1:13:A:THR:H	16	0.11
(1,115)	1:9:A:VAL:HG13	1:13:A:THR:H	16	0.11
(1,115)	1:9:A:VAL:HG21	1:13:A:THR:H	16	0.11
(1,115)	1:9:A:VAL:HG22	1:13:A:THR:H	16	0.11
(1,115)	1:9:A:VAL:HG23	1:13:A:THR:H	16	0.11
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	9	0.11
(1,84)	1:30:A:PHE:HA	1:31:A:ASP:H	19	0.11
(1,55)	1:3:A:CYS:HB2	1:4:A:TRP:H	2	0.11
(1,55)	1:3:A:CYS:HB2	1:4:A:TRP:H	6	0.11
(1,55)	1:3:A:CYS:HB2	1:4:A:TRP:H	10	0.11
(3,5)	1:13:A:THR:O	1:17:A:LEU:H	6	0.1
(1,283)	1:18:A:PRO:HG2	1:22:A:GLN:HE21	17	0.1
(1,283)	1:18:A:PRO:HG2	1:22:A:GLN:HE22	17	0.1
(1,274)	1:17:A:LEU:HD11	1:22:A:GLN:HE21	12	0.1
(1,274)	1:17:A:LEU:HD11	1:22:A:GLN:HE22	12	0.1

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,274)	1:17:A:LEU:HD12	1:22:A:GLN:HE21	12	0.1
(1,274)	1:17:A:LEU:HD12	1:22:A:GLN:HE22	12	0.1
(1,274)	1:17:A:LEU:HD13	1:22:A:GLN:HE21	12	0.1
(1,274)	1:17:A:LEU:HD13	1:22:A:GLN:HE22	12	0.1
(1,274)	1:17:A:LEU:HD21	1:22:A:GLN:HE21	12	0.1
(1,274)	1:17:A:LEU:HD21	1:22:A:GLN:HE22	12	0.1
(1,274)	1:17:A:LEU:HD22	1:22:A:GLN:HE21	12	0.1
(1,274)	1:17:A:LEU:HD22	1:22:A:GLN:HE22	12	0.1
(1,274)	1:17:A:LEU:HD23	1:22:A:GLN:HE21	12	0.1
(1,274)	1:17:A:LEU:HD23	1:22:A:GLN:HE22	12	0.1
(1,196)	1:9:A:VAL:HB	1:10:A:PRO:HD3	8	0.1
(1,196)	1:9:A:VAL:HB	1:10:A:PRO:HD3	9	0.1
(1,196)	1:9:A:VAL:HB	1:10:A:PRO:HD3	17	0.1
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	1	0.1
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	4	0.1
(1,128)	1:17:A:LEU:HB2	1:18:A:PRO:HD3	18	0.1
(1,55)	1:3:A:CYS:HB2	1:4:A:TRP:H	9	0.1
(1,32)	1:13:A:THR:HB	1:14:A:CYS:H	13	0.1

10 Dihedral-angle violation analysis [i](#)

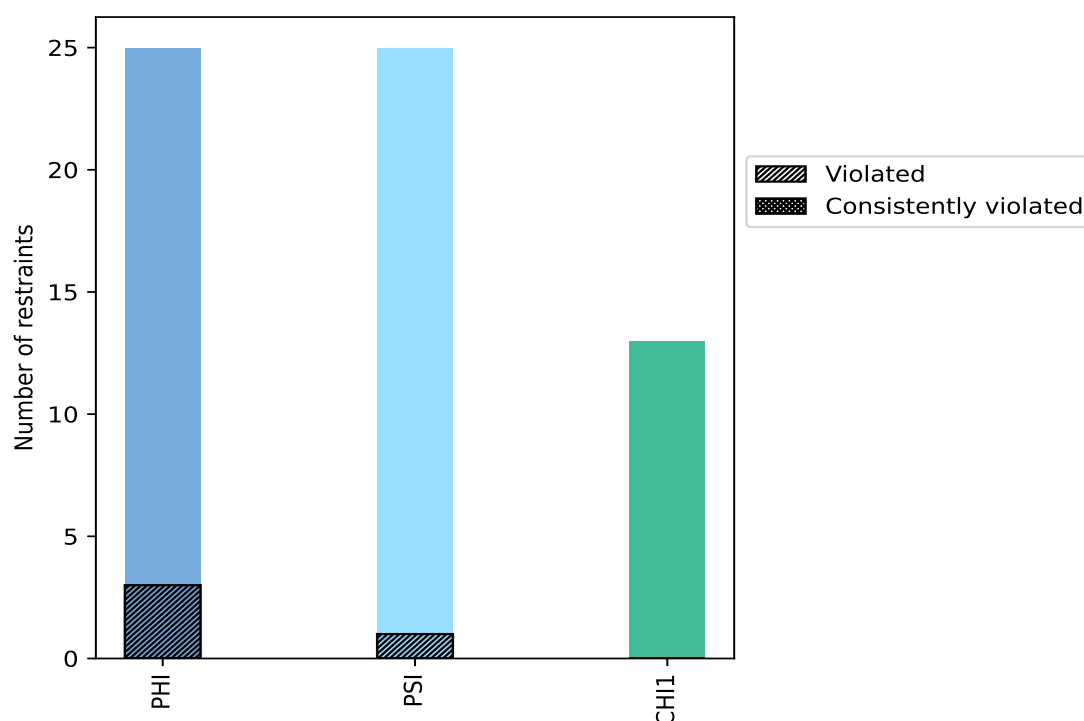
10.1 Summary of dihedral-angle violations [i](#)

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	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
PHI	25	39.7	3	12.0	4.8	0	0.0	0.0
PSI	25	39.7	1	4.0	1.6	0	0.0	0.0
CHI1	13	20.6	0	0.0	0.0	0	0.0	0.0
Total	63	100.0	4	6.3	6.3	0	0.0	0.0

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

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



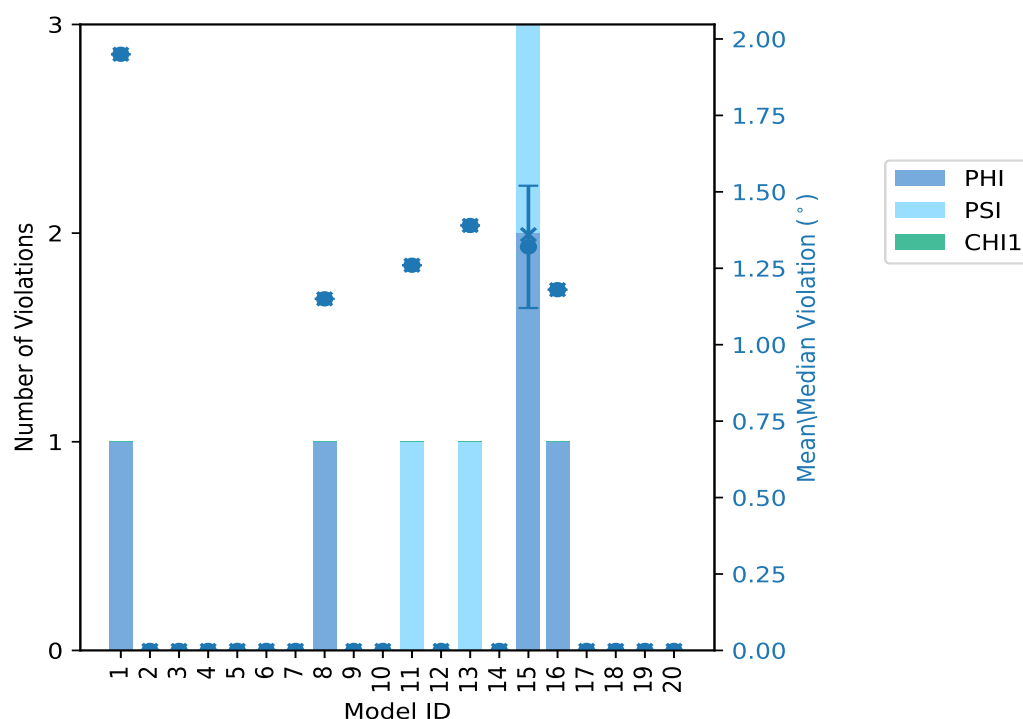
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

10.2 Dihedral-angle violation statistics for each model

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	PSI	CHI1	Total				
1	1	0	0	1	1.95	1.95	0.0	1.95
2	0	0	0	0	0.0	0.0	0.0	0.0
3	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
5	0	0	0	0	0.0	0.0	0.0	0.0
6	0	0	0	0	0.0	0.0	0.0	0.0
7	0	0	0	0	0.0	0.0	0.0	0.0
8	1	0	0	1	1.15	1.15	0.0	1.15
9	0	0	0	0	0.0	0.0	0.0	0.0
10	0	0	0	0	0.0	0.0	0.0	0.0
11	0	1	0	1	1.26	1.26	0.0	1.26
12	0	0	0	0	0.0	0.0	0.0	0.0
13	0	1	0	1	1.39	1.39	0.0	1.39
14	0	0	0	0	0.0	0.0	0.0	0.0
15	2	1	0	3	1.32	1.54	0.2	1.36
16	1	0	0	1	1.18	1.18	0.0	1.18
17	0	0	0	0	0.0	0.0	0.0	0.0
18	0	0	0	0	0.0	0.0	0.0	0.0
19	0	0	0	0	0.0	0.0	0.0	0.0
20	0	0	0	0	0.0	0.0	0.0	0.0

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	PSI	CHI1	Total	Count ¹	%
1	0	0	1	1	5.0
2	0	0	2	2	10.0
0	1	0	1	3	15.0
0	0	0	0	4	20.0
0	0	0	0	5	25.0
0	0	0	0	6	30.0
0	0	0	0	7	35.0
0	0	0	0	8	40.0
0	0	0	0	9	45.0
0	0	0	0	10	50.0
0	0	0	0	11	55.0

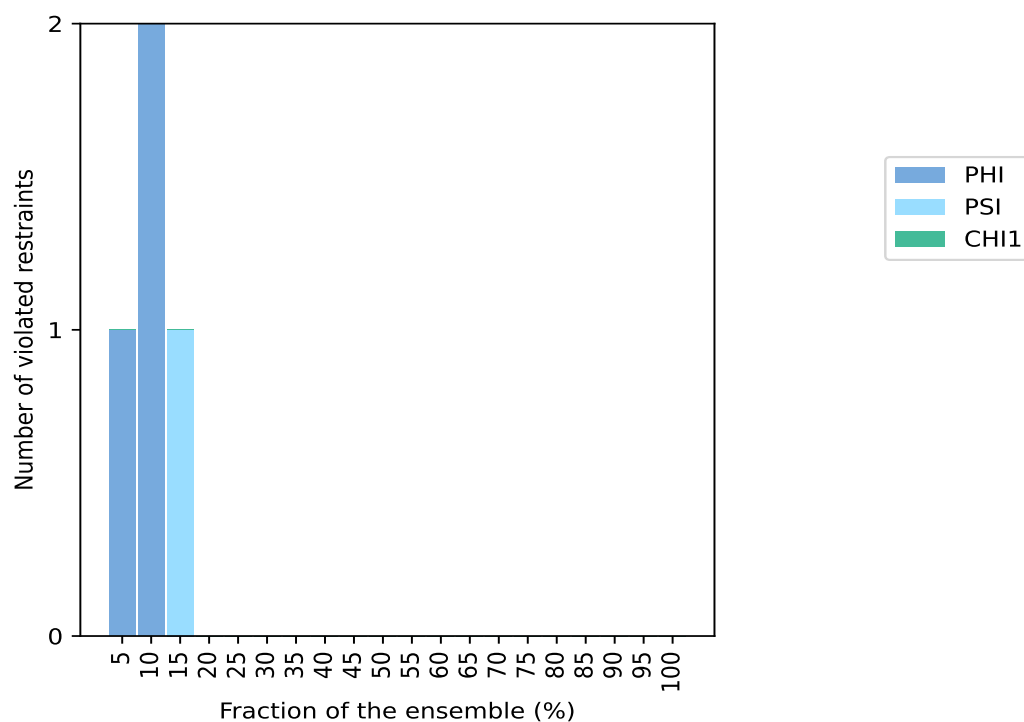
Continued on next page...

Continued from previous page...

Number of violated restraints				Fraction of the ensemble	
PHI	PSI	CHI1	Total	Count ¹	%
0	0	0	0	12	60.0
0	0	0	0	13	65.0
0	0	0	0	14	70.0
0	0	0	0	15	75.0
0	0	0	0	16	80.0
0	0	0	0	17	85.0
0	0	0	0	18	90.0
0	0	0	0	19	95.0
0	0	0	0	20	100.0

¹ Number of models with violations

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

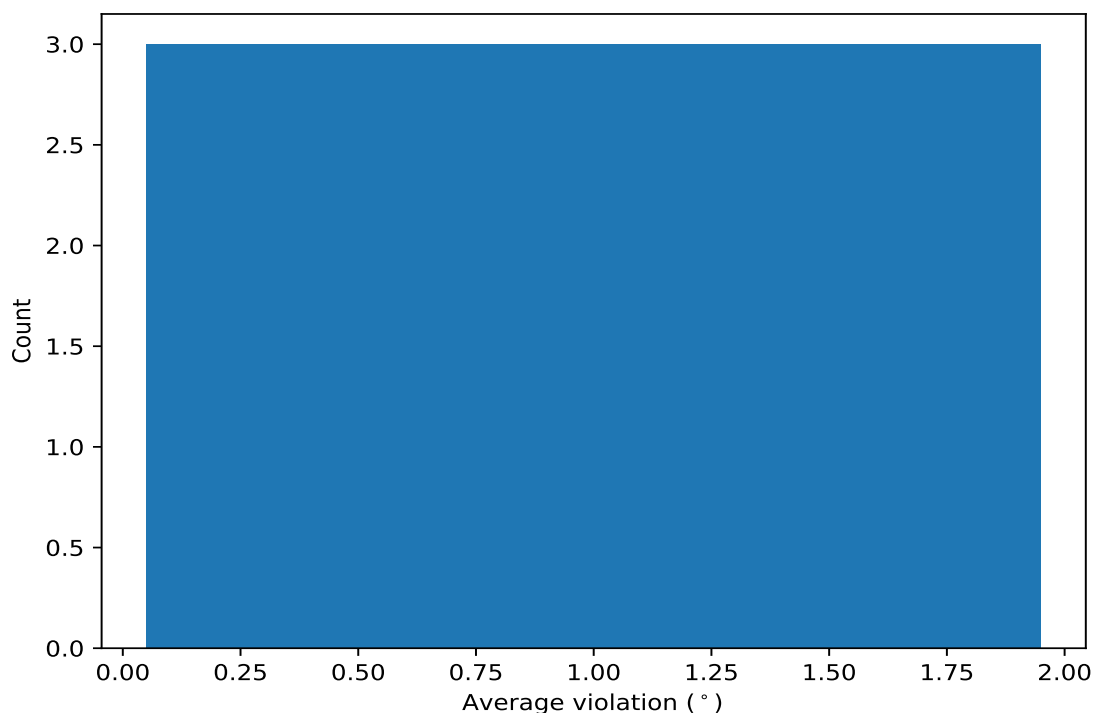


10.4 Most violated dihedral-angle restraints in the ensemble [i](#)

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.

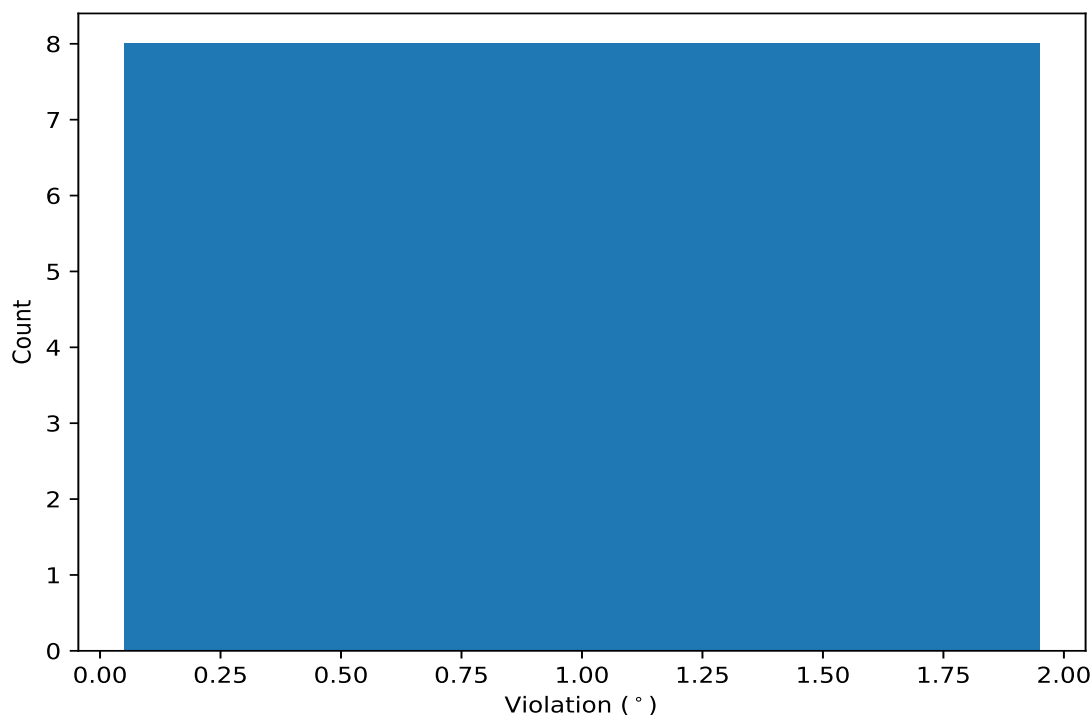
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models ¹	Mean	SD ²	Median
(1,14)	1:10:A:PRO:N	1:10:A:PRO:CA	1:10:A:PRO:C	1:11:A:TYR:N	3	1.34	0.06	1.36
(1,11)	1:8:A:CYS:C	1:9:A:VAL:N	1:9:A:VAL:CA	1:9:A:VAL:C	2	1.74	0.2	1.74
(1,15)	1:12:A:GLY:C	1:13:A:THR:N	1:13:A:THR:CA	1:13:A:THR:C	2	1.1	0.04	1.1

¹ Number of violated models, ²Standard 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 (°)
(1,11)	1:8:A:CYS:C	1:9:A:VAL:N	1:9:A:VAL:CA	1:9:A:VAL:C	1	1.95
(1,11)	1:8:A:CYS:C	1:9:A:VAL:N	1:9:A:VAL:CA	1:9:A:VAL:C	15	1.54
(1,14)	1:10:A:PRO:N	1:10:A:PRO:CA	1:10:A:PRO:C	1:11:A:TYR:N	13	1.39
(1,14)	1:10:A:PRO:N	1:10:A:PRO:CA	1:10:A:PRO:C	1:11:A:TYR:N	15	1.36
(1,14)	1:10:A:PRO:N	1:10:A:PRO:CA	1:10:A:PRO:C	1:11:A:TYR:N	11	1.26
(1,25)	1:17:A:LEU:C	1:18:A:PRO:N	1:18:A:PRO:CA	1:18:A:PRO:C	16	1.18
(1,15)	1:12:A:GLY:C	1:13:A:THR:N	1:13:A:THR:CA	1:13:A:THR:C	8	1.15
(1,15)	1:12:A:GLY:C	1:13:A:THR:N	1:13:A:THR:CA	1:13:A:THR:C	15	1.06