



wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 2MBG / pdb_00002mbg
BMRB ID : 17608
Title : Rlip76 (gap-gbd)
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Deposited on : 2013-07-30

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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 Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
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) : 2.49

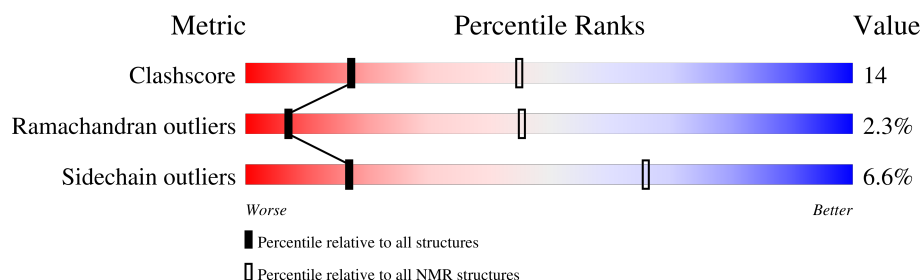
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 86%.

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	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

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	265	

2 Ensemble composition and analysis

This entry contains 35 models. Model 1 is the overall representative, medoid model (most similar to other models).

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:189-A:380, A:390-A:446 (249)	0.89	1

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 4 clusters and 1 single-model cluster was found.

Cluster number	Models
1	1, 2, 3, 4, 5, 6, 9, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20, 24, 25, 27, 28, 29, 30, 31, 32, 34
2	7, 23, 26, 35
3	8, 21
4	22, 33
Single-model clusters	10

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 4406 atoms, of which 2234 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called RalA-binding protein 1.

Mol	Chain	Residues	Atoms						Trace
1	A	265	Total	C	H	N	O	S	0
			4406	1385	2234	373	399	15	

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	182	HIS	-	expression tag	UNP Q15311
A	183	MET	-	expression tag	UNP Q15311

5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 35 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
CNS	refinement	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	3268
Number of shifts mapped to atoms	3268
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	86%

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	0.43±0.01	0±0/2074 (0.0± 0.0%)	0.75±0.01	0±0/2795 (0.0± 0.0%)
All	All	0.43	0/72590 (0.0%)	0.75	3/97825 (0.0%)

There are no bond-length outliers.

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	368	PHE	CB-CA-C	-5.52	107.50	114.40	9	3

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	2041	2100	2096	56±5
All	All	71435	73500	73360	1976

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 14.

5 of 589 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:247:TYR:HA	1:A:251:GLU:HB3	0.90	1.44	32	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:241:ASP:HA	1:A:244:LYS:HE2	0.89	1.43	14	5
1:A:295:THR:HB	1:A:298:GLU:HG2	0.88	1.46	12	22
1:A:225:MET:HG3	1:A:331:LYS:HG3	0.86	1.45	14	11
1:A:300:VAL:HG12	1:A:363:HIS:HB2	0.86	1.48	17	13

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	248/265 (94%)	226±3 (91±1%)	16±3 (7±1%)	6±2 (2±1%)	7	45
All	All	8680/9275 (94%)	7914 (91%)	566 (7%)	200 (2%)	7	45

5 of 24 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	207	ILE	33
1	A	280	THR	31
1	A	390	THR	25
1	A	204	TYR	22
1	A	189	ILE	21

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	228/243 (94%)	213±2 (93±1%)	15±2 (7±1%)	17	66
All	All	7980/8505 (94%)	7455 (93%)	525 (7%)	17	66

5 of 85 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	279	LEU	35
1	A	340	GLN	35
1	A	306	LEU	34
1	A	370	ASN	24
1	A	364	VAL	22

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 86% for the well-defined parts and 85% 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	3268
Number of shifts mapped to atoms	3268
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	8

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	254	-0.50 ± 0.09	Should be checked
$^{13}\text{C}_\beta$	242	0.23 ± 0.06	None needed (< 0.5 ppm)
$^{13}\text{C}'$	232	-0.36 ± 0.09	None needed (< 0.5 ppm)
^{15}N	246	0.96 ± 0.20	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 86%, i.e. 3106 atoms were assigned a chemical shift out of a possible 3624. 0 out of 51 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	1195/1237 (97%)	490/499 (98%)	467/498 (94%)	238/240 (99%)
Sidechain	1748/2170 (81%)	1199/1407 (85%)	531/674 (79%)	18/89 (20%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	163/217 (75%)	88/104 (85%)	73/103 (71%)	2/10 (20%)
Overall	3106/3624 (86%)	1777/2010 (88%)	1071/1275 (84%)	258/339 (76%)

7.1.4 Statistically unusual chemical shifts [i](#)

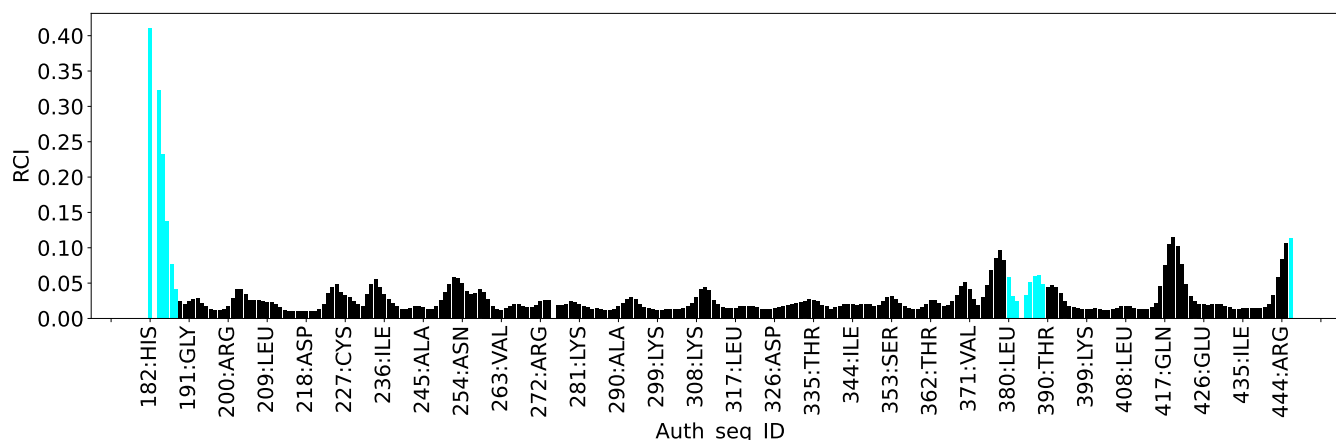
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	271	LEU	CG	66.14	21.37 – 32.19	36.4
1	A	280	THR	HG1	6.12	0.08 – 2.19	23.6
1	A	201	THR	HG1	4.45	0.08 – 2.19	15.7
1	A	335	THR	HG1	3.57	0.08 – 2.19	11.5
1	A	244	LYS	HG2	-0.54	0.13 – 2.61	-7.7
1	A	209	LEU	HD11	-0.90	-0.61 – 2.12	-6.1
1	A	209	LEU	HD12	-0.90	-0.61 – 2.12	-6.1
1	A	209	LEU	HD13	-0.90	-0.61 – 2.12	-6.1

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	7318
Intra-residue ($ i-j =0$)	2430
Sequential ($ i-j =1$)	1662
Medium range ($ i-j >1$ and $ i-j <5$)	1458
Long range ($ i-j \geq 5$)	1548
Inter-chain	0
Hydrogen bond restraints	220
Disulfide bond restraints	0
Total dihedral-angle restraints	434
Number of unmapped restraints	0
Number of restraints per residue	29.3
Number of long range restraints per residue ¹	5.8

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)	94.4	0.2
0.2-0.5 (Medium)	135.3	0.5
>0.5 (Large)	210.9	5.14

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)	41.7	9.88
10.0-20.0 (Medium)	0.7	17.97
>20.0 (Large)	None	None

9 Distance violation analysis ⓘ

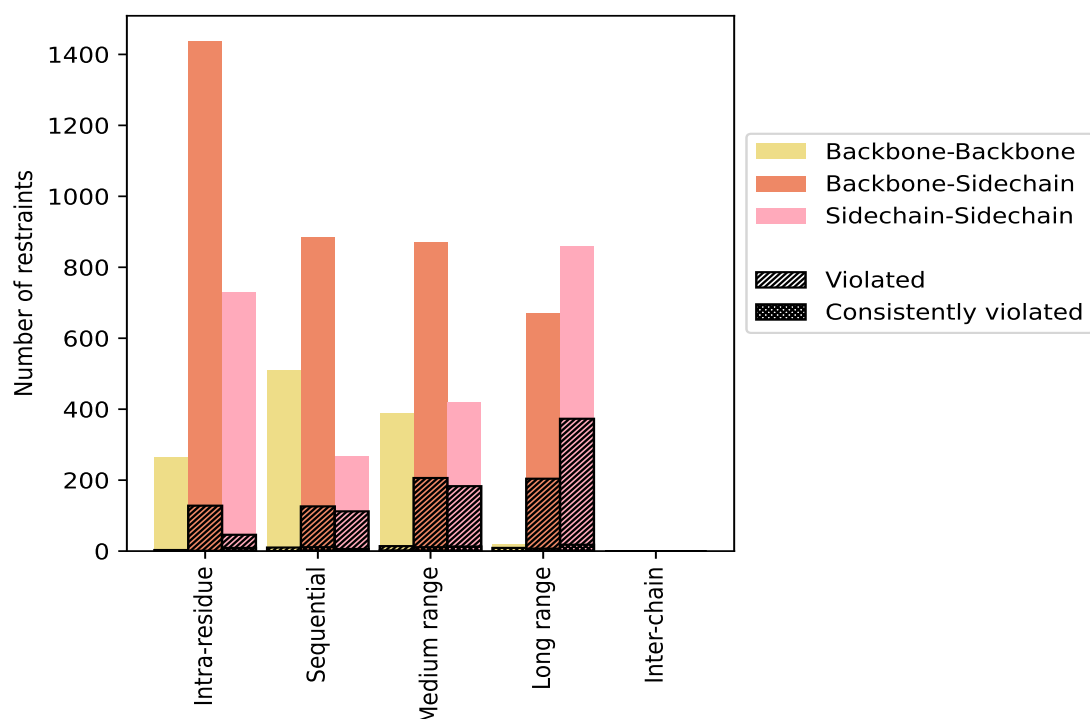
9.1 Summary of distance violations ⓘ

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.

Restraints type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	2430	33.2	177	7.3	2.4	10	0.4	0.1
Backbone-Backbone	265	3.6	3	1.1	0.0	0	0.0	0.0
Backbone-Sidechain	1437	19.6	128	8.9	1.7	1	0.1	0.0
Sidechain-Sidechain	728	9.9	46	6.3	0.6	9	1.2	0.1
Sequential ($i-j =1$)	1662	22.7	248	14.9	3.4	17	1.0	0.2
Backbone-Backbone	510	7.0	10	2.0	0.1	0	0.0	0.0
Backbone-Sidechain	885	12.1	126	14.2	1.7	11	1.2	0.2
Sidechain-Sidechain	267	3.6	112	41.9	1.5	6	2.2	0.1
Medium range ($i-j >1$ & $i-j <5$)	1458	19.9	379	26.0	5.2	23	1.6	0.3
Backbone-Backbone	389	5.3	14	3.6	0.2	2	0.5	0.0
Backbone-Sidechain	651	8.9	182	28.0	2.5	8	1.2	0.1
Sidechain-Sidechain	418	5.7	183	43.8	2.5	13	3.1	0.2
Long range ($i-j \geq 5$)	1548	21.2	586	37.9	8.0	25	1.6	0.3
Backbone-Backbone	20	0.3	9	45.0	0.1	0	0.0	0.0
Backbone-Sidechain	669	9.1	204	30.5	2.8	7	1.0	0.1
Sidechain-Sidechain	859	11.7	373	43.4	5.1	18	2.1	0.2
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	220	3.0	24	10.9	0.3	3	1.4	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	7318	100.0	1414	19.3	19.3	78	1.1	1.1
Backbone-Backbone	1184	16.2	36	3.0	0.5	2	0.2	0.0
Backbone-Sidechain	3862	52.8	664	17.2	9.1	30	0.8	0.4
Sidechain-Sidechain	2272	31.0	714	31.4	9.8	46	2.0	0.6

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 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 ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	53	79	129	197	0	458	0.64	4.52	0.58	0.47
2	44	71	128	198	0	441	0.68	4.68	0.62	0.5
3	44	88	136	186	0	454	0.63	4.38	0.54	0.46
4	48	73	125	192	0	438	0.64	3.38	0.53	0.47
5	37	66	133	181	0	417	0.61	3.62	0.5	0.49
6	44	71	134	204	0	453	0.62	4.02	0.51	0.48
7	45	77	132	219	0	473	0.67	4.41	0.57	0.49
8	38	80	133	177	0	428	0.61	2.83	0.5	0.46
9	45	75	132	183	0	435	0.59	4.3	0.54	0.41
10	49	73	149	198	0	469	0.67	4.09	0.57	0.51

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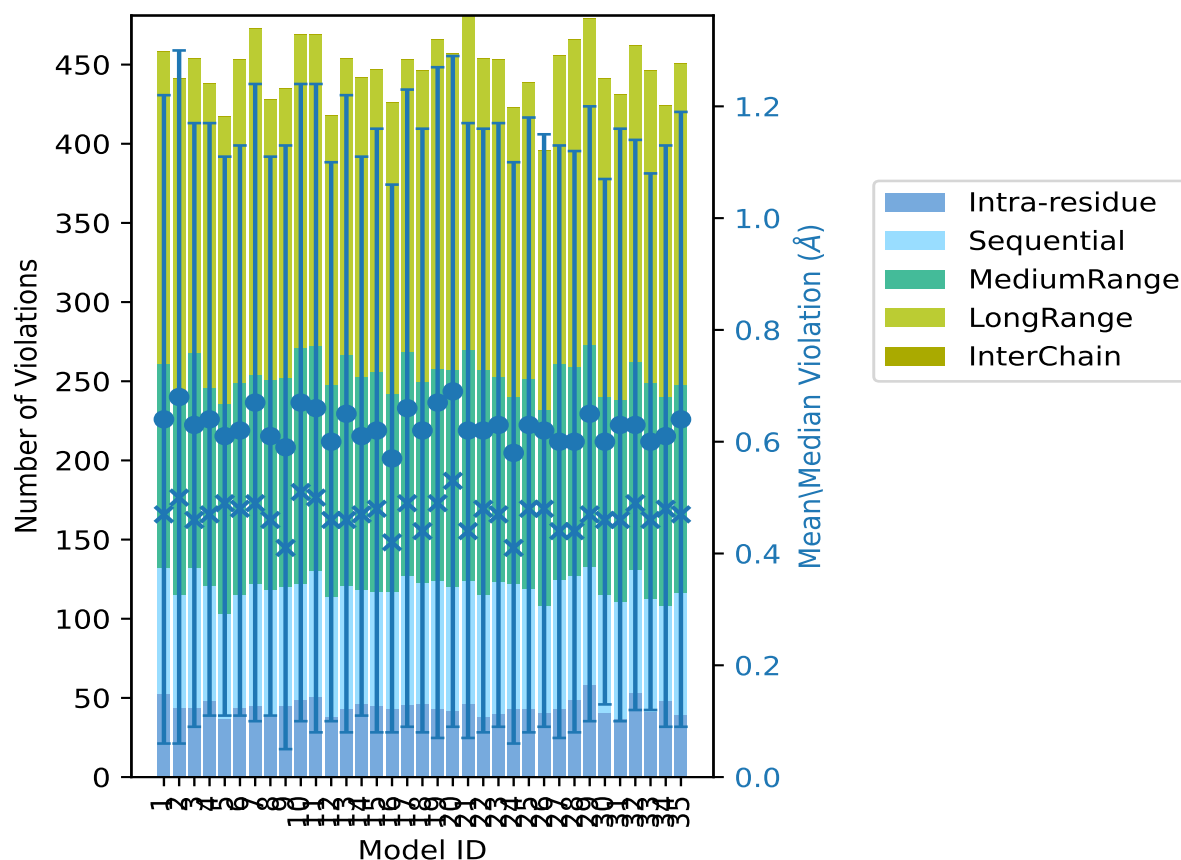
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
11	51	79	142	197	0	469	0.66	5.14	0.58	0.5
12	38	76	134	170	0	418	0.6	3.33	0.5	0.46
13	43	78	146	187	0	454	0.65	4.69	0.57	0.46
14	46	72	135	189	0	442	0.61	3.0	0.5	0.47
15	45	72	139	191	0	447	0.62	5.14	0.54	0.48
16	43	74	125	184	0	426	0.57	4.78	0.49	0.42
17	46	81	142	184	0	453	0.66	4.83	0.57	0.49
18	46	77	127	196	0	446	0.62	4.95	0.54	0.44
19	43	81	134	208	0	466	0.67	3.79	0.6	0.49
20	42	78	137	200	0	457	0.69	4.65	0.6	0.53
21	46	78	146	211	0	481	0.62	4.77	0.55	0.44
22	38	77	142	197	0	454	0.62	4.75	0.54	0.48
23	40	83	130	200	0	453	0.63	4.25	0.54	0.47
24	43	79	118	183	0	423	0.58	4.17	0.52	0.41
25	43	76	133	187	0	439	0.63	4.18	0.55	0.48
26	41	67	124	164	0	396	0.62	4.63	0.53	0.48
27	43	82	136	195	0	456	0.6	4.89	0.53	0.44
28	49	78	132	207	0	466	0.6	3.51	0.52	0.44
29	58	75	140	206	0	479	0.65	4.19	0.55	0.47
30	41	74	125	201	0	441	0.6	3.12	0.47	0.46
31	37	74	127	193	0	431	0.63	3.57	0.53	0.46
32	53	78	131	200	0	462	0.63	2.92	0.51	0.49
33	41	72	136	197	0	446	0.6	2.69	0.48	0.46
34	48	60	132	184	0	424	0.61	4.32	0.52	0.48
35	39	77	132	203	0	451	0.64	4.42	0.55	0.47

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 [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, 5708(IR:2253, SQ:1414, MR:1079, LR:962, 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	%
41	61	62	108	0	272	1	2.9
18	19	43	56	0	136	2	5.7
16	18	25	40	0	99	3	8.6
13	14	17	40	0	84	4	11.4
8	9	15	23	0	55	5	14.3
6	8	16	24	0	54	6	17.1

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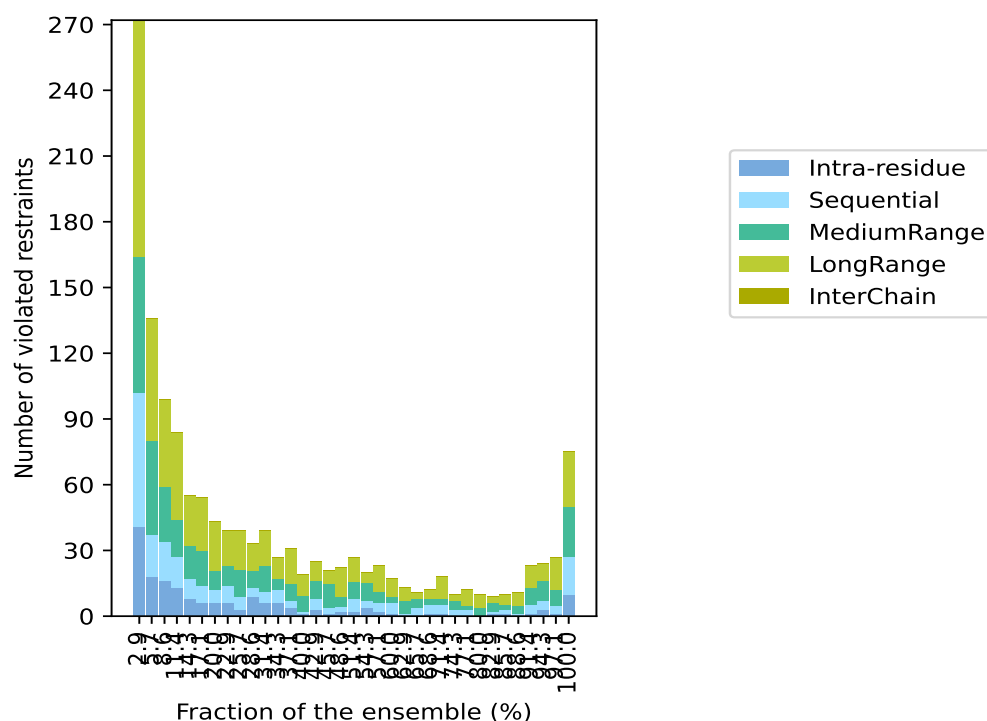
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Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
6	6	9	22	0	43	7	20.0
6	8	9	16	0	39	8	22.9
3	6	12	18	0	39	9	25.7
9	4	8	12	0	33	10	28.6
6	5	12	16	0	39	11	31.4
6	6	5	10	0	27	12	34.3
4	3	8	16	0	31	13	37.1
1	1	7	10	0	19	14	40.0
3	5	8	9	0	25	15	42.9
1	3	11	6	0	21	16	45.7
2	2	5	13	0	22	17	48.6
2	6	8	11	0	27	18	51.4
4	3	8	5	0	20	19	54.3
2	4	5	12	0	23	20	57.1
1	5	3	8	0	17	21	60.0
0	1	6	6	0	13	22	62.9
0	4	4	3	0	11	23	65.7
1	4	3	4	0	12	24	68.6
1	4	3	10	0	18	25	71.4
0	3	4	3	0	10	26	74.3
0	3	2	7	0	12	27	77.1
0	0	4	6	0	10	28	80.0
1	1	4	3	0	9	29	82.9
1	2	2	5	0	10	30	85.7
0	1	4	6	0	11	31	88.6
1	4	8	10	0	23	32	91.4
3	4	9	8	0	24	33	94.3
1	4	7	15	0	27	34	97.1
10	17	23	25	0	75	35	100.0

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

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

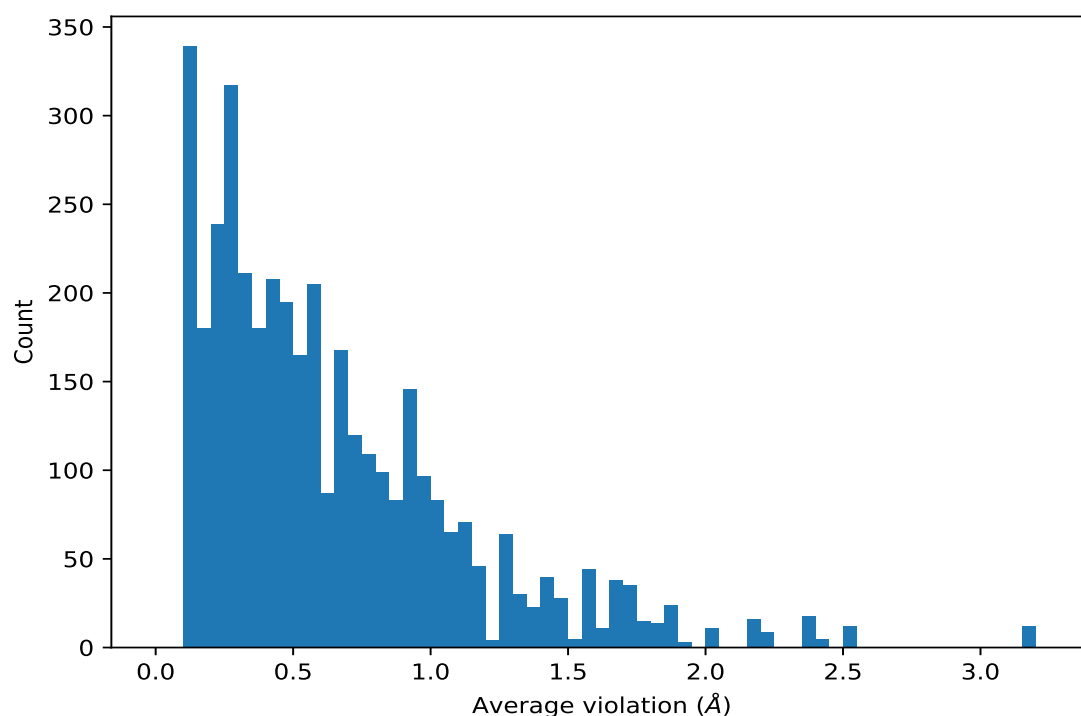
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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 violations for the 10 worst performing restraints, sorted by number of violated models and the mean violation 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)
(1,5791)	1:367:A:LEU:HD22	1:318:A:ILE:HG21	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD22	1:318:A:ILE:HG22	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD23	1:318:A:ILE:HG22	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD22	1:318:A:ILE:HG23	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD23	1:318:A:ILE:HG21	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD21	1:318:A:ILE:HG21	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD21	1:318:A:ILE:HG22	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD21	1:323:A:VAL:HG22	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD22	1:323:A:VAL:HG21	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD23	1:323:A:VAL:HG21	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD21	1:318:A:ILE:HG23	35	3.16	1.45	4.09
(1,5791)	1:367:A:LEU:HD23	1:318:A:ILE:HG23	35	3.16	1.45	4.09
(1,5949)	1:407:A:PHE:HD2	1:436:A:LEU:HD21	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD2	1:408:A:LEU:HD21	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD2	1:436:A:LEU:HD23	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD1	1:436:A:LEU:HD23	35	2.54	0.43	2.55

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,5949)	1:407:A:PHE:HD2	1:408:A:LEU:HD22	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD2	1:436:A:LEU:HD22	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD1	1:412:A:LEU:HD13	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD2	1:408:A:LEU:HD23	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD1	1:408:A:LEU:HD23	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD1	1:436:A:LEU:HD21	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD1	1:412:A:LEU:HD12	35	2.54	0.43	2.55
(1,5949)	1:407:A:PHE:HD1	1:436:A:LEU:HD22	35	2.54	0.43	2.55
(1,6499)	1:271:A:LEU:HD12	1:320:A:TRP:HB2	35	2.39	1.01	2.07
(1,6499)	1:271:A:LEU:HD11	1:210:A:PRO:HD3	35	2.39	1.01	2.07
(1,6499)	1:271:A:LEU:HD13	1:210:A:PRO:HD3	35	2.39	1.01	2.07
(1,6499)	1:271:A:LEU:HD12	1:210:A:PRO:HD3	35	2.39	1.01	2.07
(1,6499)	1:271:A:LEU:HD11	1:320:A:TRP:HB2	35	2.39	1.01	2.07
(1,6499)	1:271:A:LEU:HD13	1:320:A:TRP:HB2	35	2.39	1.01	2.07
(1,6012)	1:359:A:VAL:HG23	1:364:A:VAL:HG12	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG22	1:364:A:VAL:HG13	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG21	1:364:A:VAL:HG12	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG23	1:364:A:VAL:HG13	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG23	1:364:A:VAL:HG11	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG21	1:367:A:LEU:HB2	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG21	1:364:A:VAL:HG13	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG22	1:364:A:VAL:HG11	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG22	1:367:A:LEU:HB2	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG22	1:364:A:VAL:HG12	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG21	1:364:A:VAL:HG11	35	2.38	0.82	2.13
(1,6012)	1:359:A:VAL:HG23	1:367:A:LEU:HB2	35	2.38	0.82	2.13
(1,6178)	1:364:A:VAL:HG21	1:279:A:LEU:HD11	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG23	1:279:A:LEU:HD11	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG22	1:279:A:LEU:HD11	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG23	1:279:A:LEU:HD13	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG22	1:279:A:LEU:HD13	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG23	1:279:A:LEU:HD12	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG21	1:300:A:VAL:HG13	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG21	1:279:A:LEU:HD13	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG21	1:279:A:LEU:HD12	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG21	1:300:A:VAL:HG12	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG22	1:300:A:VAL:HG11	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG22	1:300:A:VAL:HG12	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG22	1:300:A:VAL:HG13	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG23	1:300:A:VAL:HG11	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG23	1:300:A:VAL:HG12	35	2.19	0.32	2.1
(1,6178)	1:364:A:VAL:HG22	1:279:A:LEU:HD12	35	2.19	0.32	2.1

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,6718)	1:359:A:VAL:HG23	1:322:A:ILE:HG22	35	2.0	0.65	1.84
(1,6718)	1:359:A:VAL:HG23	1:322:A:ILE:HG21	35	2.0	0.65	1.84
(1,6718)	1:359:A:VAL:HG21	1:322:A:ILE:HG21	35	2.0	0.65	1.84
(1,6718)	1:359:A:VAL:HG21	1:322:A:ILE:HG22	35	2.0	0.65	1.84
(1,6718)	1:359:A:VAL:HG23	1:322:A:ILE:HG23	35	2.0	0.65	1.84
(1,6718)	1:359:A:VAL:HG22	1:322:A:ILE:HG21	35	2.0	0.65	1.84
(1,6718)	1:359:A:VAL:HG21	1:322:A:ILE:HG23	35	2.0	0.65	1.84
(1,6718)	1:359:A:VAL:HG22	1:322:A:ILE:HG22	35	2.0	0.65	1.84
(1,6056)	1:373:A:LEU:HD21	1:325:A:MET:HE1	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD22	1:325:A:MET:HE1	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD22	1:321:A:LEU:HD11	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD22	1:325:A:MET:HE3	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD23	1:325:A:MET:HE1	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD22	1:321:A:LEU:HD12	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD23	1:325:A:MET:HE2	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD23	1:325:A:MET:HE3	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD22	1:325:A:MET:HE2	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD21	1:364:A:VAL:HG23	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD23	1:364:A:VAL:HG23	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD21	1:325:A:MET:HE2	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD22	1:207:A:ILE:HD13	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD23	1:364:A:VAL:HG22	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD23	1:364:A:VAL:HG21	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD22	1:364:A:VAL:HG22	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD22	1:364:A:VAL:HG23	35	1.87	0.42	1.9
(1,6056)	1:373:A:LEU:HD21	1:325:A:MET:HE3	35	1.87	0.42	1.9
(1,4897)	1:313:A:CYS:H	1:209:A:LEU:HD11	35	1.83	0.89	1.49
(1,4897)	1:313:A:CYS:H	1:209:A:LEU:HD12	35	1.83	0.89	1.49
(1,4897)	1:313:A:CYS:H	1:317:A:LEU:HD12	35	1.83	0.89	1.49
(1,4897)	1:313:A:CYS:H	1:317:A:LEU:HD11	35	1.83	0.89	1.49
(1,4897)	1:313:A:CYS:H	1:317:A:LEU:HD13	35	1.83	0.89	1.49
(1,4897)	1:313:A:CYS:H	1:209:A:LEU:HD13	35	1.83	0.89	1.49
(1,5723)	1:203:A:MET:HE2	1:217:A:ILE:HG21	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE1	1:217:A:ILE:HG21	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE3	1:217:A:ILE:HG12	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE2	1:217:A:ILE:HG22	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE3	1:217:A:ILE:HG23	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE2	1:217:A:ILE:HG23	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE2	1:207:A:ILE:HG23	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE1	1:217:A:ILE:HG23	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE3	1:217:A:ILE:HG22	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE3	1:217:A:ILE:HG21	35	1.72	0.23	1.72

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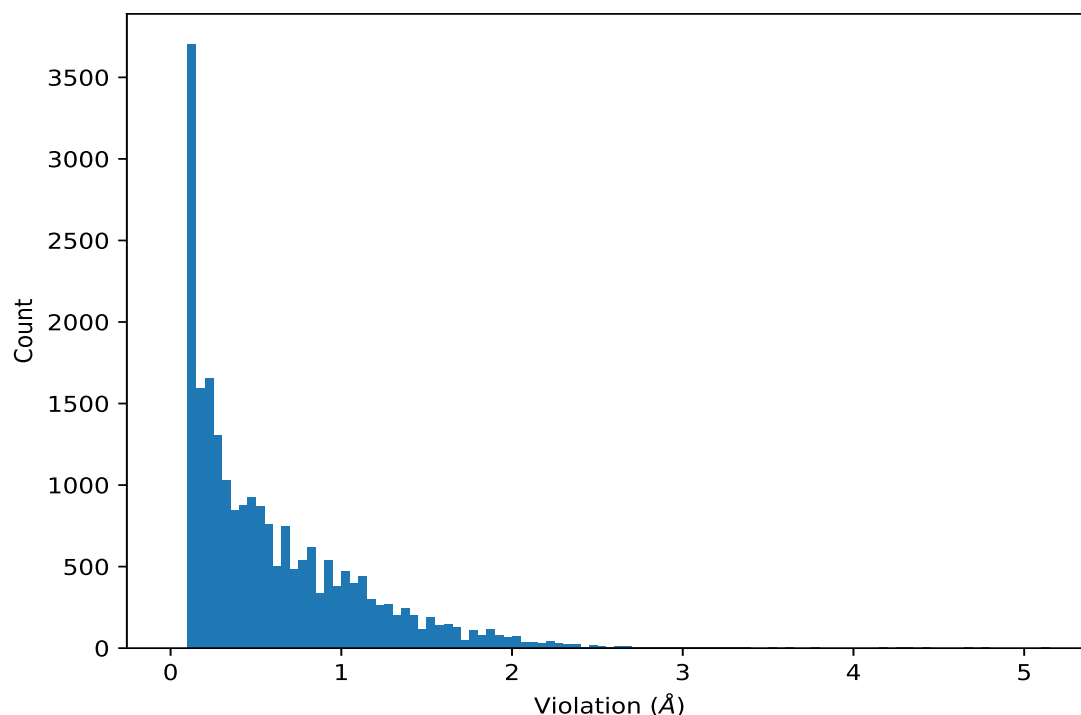
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,5723)	1:203:A:MET:HE1	1:217:A:ILE:HG22	35	1.72	0.23	1.72
(1,5723)	1:203:A:MET:HE2	1:207:A:ILE:HG22	35	1.72	0.23	1.72
(1,5756)	1:203:A:MET:HE2	1:255:A:LEU:HD13	35	1.66	0.63	1.74

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 provides the 10 worst performing restraints, sorted by the violation 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)
(1,6499)	1:271:A:LEU:HD13	1:210:A:PRO:HD3	11	5.14
(1,6499)	1:271:A:LEU:HD11	1:210:A:PRO:HD3	15	5.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,5791)	1:367:A:LEU:HD21	1:318:A:ILE:HG21	18	4.95
(1,5791)	1:367:A:LEU:HD23	1:318:A:ILE:HG22	27	4.89
(1,5791)	1:367:A:LEU:HD22	1:318:A:ILE:HG21	17	4.83
(1,6499)	1:271:A:LEU:HD12	1:210:A:PRO:HD3	16	4.78
(1,5791)	1:367:A:LEU:HD22	1:323:A:VAL:HG21	21	4.77
(1,5791)	1:367:A:LEU:HD23	1:323:A:VAL:HG21	22	4.75
(1,5791)	1:367:A:LEU:HD22	1:318:A:ILE:HG21	13	4.69
(1,6012)	1:359:A:VAL:HG23	1:364:A:VAL:HG12	2	4.68

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

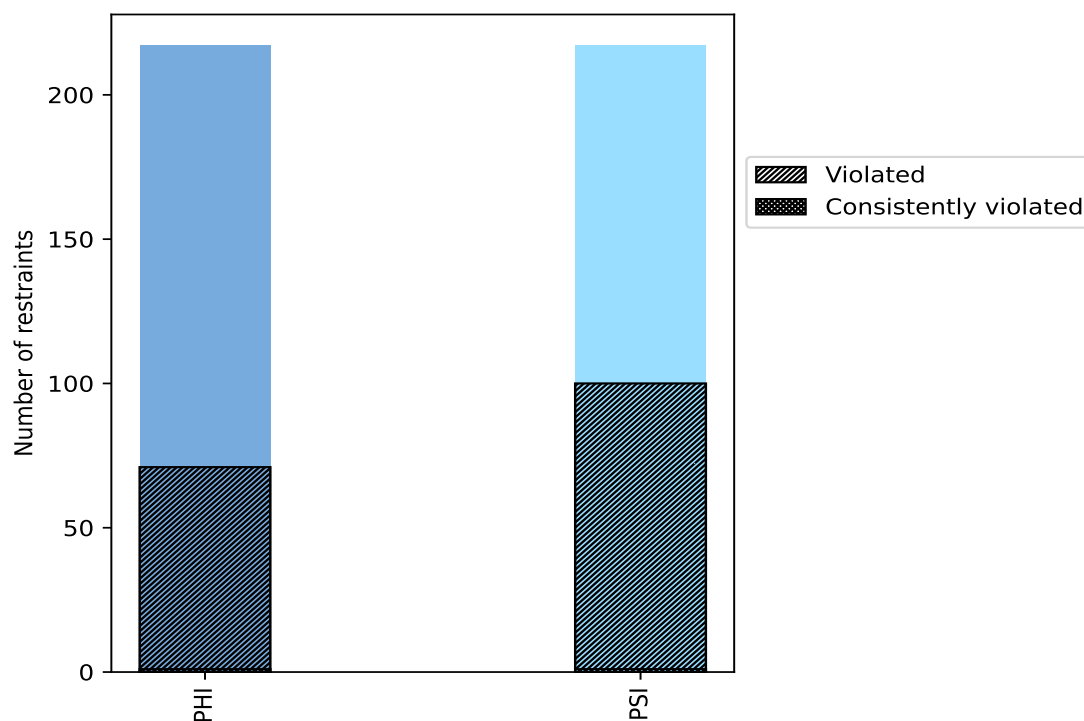
10.1 Summary of dihedral-angle violations [i](#)

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	217	50.0	71	32.7	16.4	1	0.5	0.2
PSI	217	50.0	100	46.1	23.0	1	0.5	0.2
Total	434	100.0	171	39.4	39.4	2	0.5	0.5

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](#)



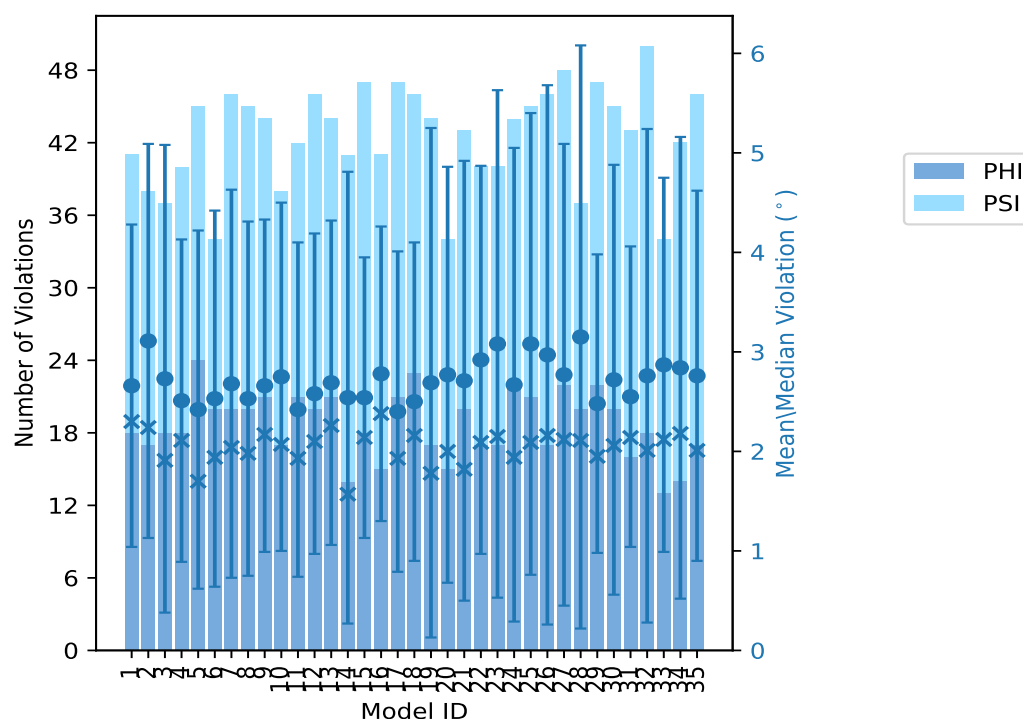
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

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	Total				
1	18	23	41	2.66	9.47	1.62	2.3
2	17	21	38	3.11	8.68	1.98	2.24
3	18	19	37	2.73	12.29	2.35	1.91
4	18	22	40	2.51	10.09	1.62	2.11
5	24	21	45	2.42	10.68	1.8	1.7
6	20	14	34	2.53	10.41	1.89	1.94
7	20	26	46	2.68	12.39	1.95	2.04
8	20	25	45	2.53	9.73	1.78	1.98
9	21	23	44	2.66	8.4	1.67	2.17
10	17	21	38	2.75	9.57	1.75	2.07
11	21	21	42	2.42	9.88	1.68	1.93
12	20	26	46	2.58	8.65	1.61	2.1
13	21	23	44	2.69	9.63	1.63	2.26
14	14	27	41	2.54	11.11	2.27	1.57
15	18	29	47	2.54	8.48	1.41	2.14
16	15	26	41	2.78	7.68	1.48	2.38
17	21	26	47	2.4	10.59	1.61	1.93
18	23	23	46	2.5	9.81	1.6	2.16
19	17	27	44	2.69	14.43	2.56	1.78
20	15	19	34	2.77	11.85	2.09	2.0
21	20	23	43	2.71	11.19	2.21	1.82
22	17	23	40	2.92	8.65	1.95	2.09
23	17	23	40	3.08	13.99	2.55	2.15
24	22	22	44	2.67	14.26	2.38	1.94
25	21	24	45	3.08	11.98	2.32	2.09
26	17	29	46	2.97	17.97	2.71	2.16
27	22	26	48	2.77	12.0	2.32	2.12
28	20	17	37	3.15	17.0	2.93	2.11
29	22	25	47	2.48	8.2	1.5	1.95
30	20	25	45	2.72	12.83	2.16	2.06
31	16	27	43	2.55	9.31	1.51	2.14
32	18	32	50	2.76	14.8	2.48	2.01
33	13	21	34	2.87	8.39	1.88	2.12
34	14	28	42	2.84	14.18	2.32	2.18
35	17	29	46	2.76	9.49	1.86	2.01

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	Total	Count ¹	%
19	22	41	1	2.9
7	14	21	2	5.7
8	10	18	3	8.6
4	2	6	4	11.4
1	5	6	5	14.3
1	4	5	6	17.1
5	4	9	7	20.0
1	4	5	8	22.9
2	5	7	9	25.7
0	1	1	10	28.6
1	4	5	11	31.4

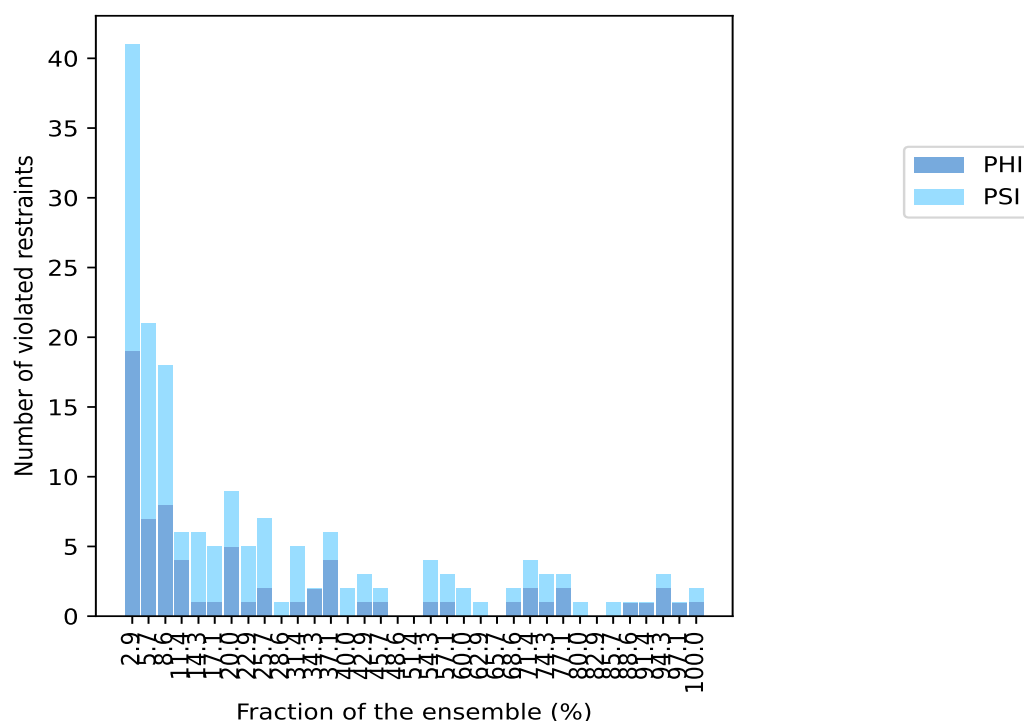
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Number of violated restraints			Fraction of the ensemble	
PHI	PSI	Total	Count ¹	%
2	0	2	12	34.3
4	2	6	13	37.1
0	2	2	14	40.0
1	2	3	15	42.9
1	1	2	16	45.7
0	0	0	17	48.6
0	0	0	18	51.4
1	3	4	19	54.3
1	2	3	20	57.1
0	2	2	21	60.0
0	1	1	22	62.9
0	0	0	23	65.7
1	1	2	24	68.6
2	2	4	25	71.4
1	2	3	26	74.3
2	1	3	27	77.1
0	1	1	28	80.0
0	0	0	29	82.9
0	1	1	30	85.7
1	0	1	31	88.6
1	0	1	32	91.4
2	1	3	33	94.3
1	0	1	34	97.1
1	1	2	35	100.0

¹ Number of models with violations

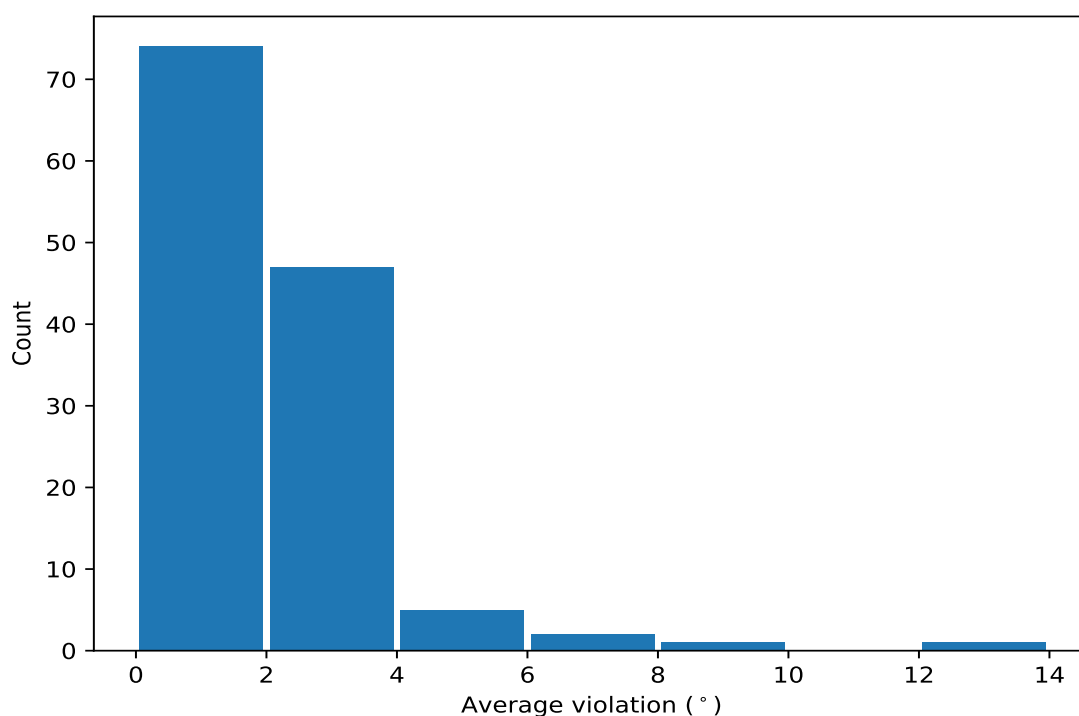
10.3.1 Bar graph : Dihedral-angle Violation statistics for the ensemble [i](#)



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 violations for the 10 worst performing restraints, sorted by number of violated models and the mean violation value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint.

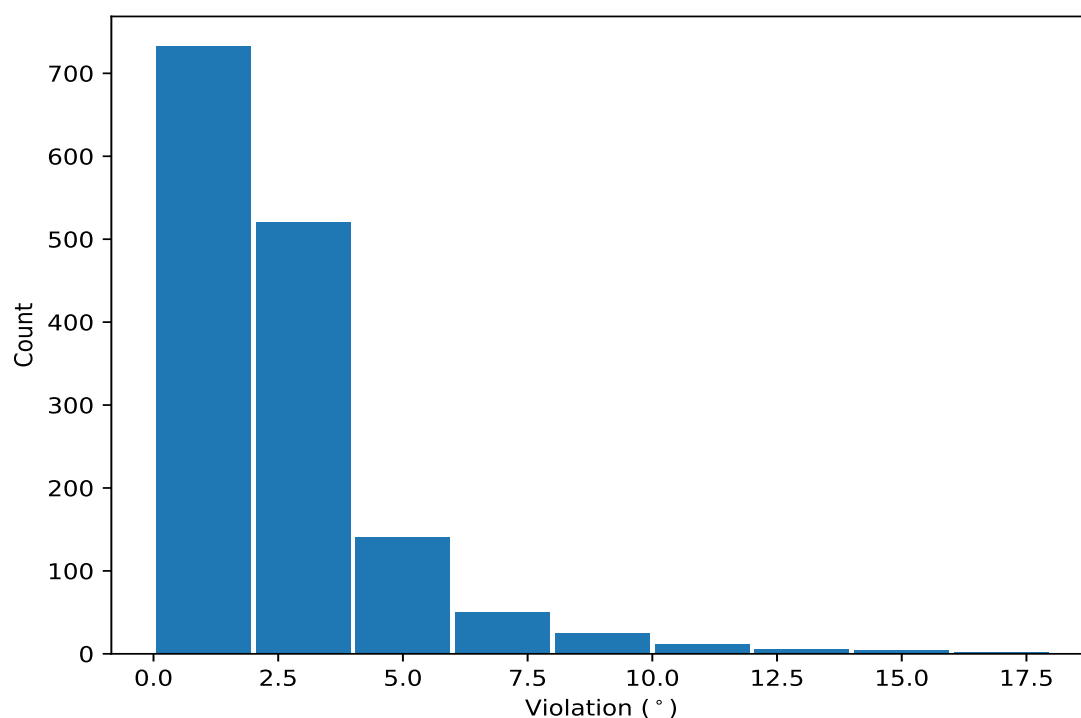
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models ¹	Mean	SD ²	Median
(1,295)	1:357:A:LEU:C	1:358:A:TYR:N	1:358:A:TYR:CA	1:358:A:TYR:C	35	9.19	1.11	9.33
(1,302)	1:361:A:PHE:N	1:361:A:PHE:CA	1:361:A:PHE:C	1:362:A:THR:N	35	5.07	1.28	5.21
(1,11)	1:195:A:ALA:C	1:196:A:ASP:N	1:196:A:ASP:CA	1:196:A:ASP:C	34	3.85	1.14	3.76
(1,135)	1:269:A:GLN:C	1:270:A:TYR:N	1:270:A:TYR:CA	1:270:A:TYR:C	33	5.42	1.13	5.47
(1,257)	1:337:A:MET:C	1:338:A:ASN:N	1:338:A:ASN:CA	1:338:A:ASN:C	33	3.7	1.15	3.55
(1,124)	1:264:A:ALA:N	1:264:A:ALA:CA	1:264:A:ALA:C	1:265:A:SER:N	33	2.74	0.94	2.46
(1,407)	1:430:A:TRP:C	1:431:A:GLU:N	1:431:A:GLU:CA	1:431:A:GLU:C	32	1.96	0.6	1.87
(1,79)	1:238:A:SER:C	1:239:A:LYS:N	1:239:A:LYS:CA	1:239:A:LYS:C	31	3.33	1.08	3.2
(1,162)	1:287:A:PHE:N	1:287:A:PHE:CA	1:287:A:PHE:C	1:288:A:GLU:N	30	2.15	0.7	2.14
(1,184)	1:299:A:LYS:N	1:299:A:LYS:CA	1:299:A:LYS:C	1:300:A:VAL:N	28	2.39	0.7	2.53

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 provides the list of violations for the 10 worst performing restraints, sorted by the violation 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 (°)
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	26	17.97
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	28	17.0
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	32	14.8
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	19	14.43
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	24	14.26
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	34	14.18
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	23	13.99
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	30	12.83
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	7	12.39
(1,256)	1:335:A:THR:N	1:335:A:THR:CA	1:335:A:THR:C	1:336:A:LYS:N	3	12.29