



wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 1HDL / pdb_00001hdl
BMRB ID : 4910
Title : LEKTI domain one
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Deposited on : 2000-11-16

This is a wwPDB NMR Structure Validation Summary 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 Phenix2.0
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

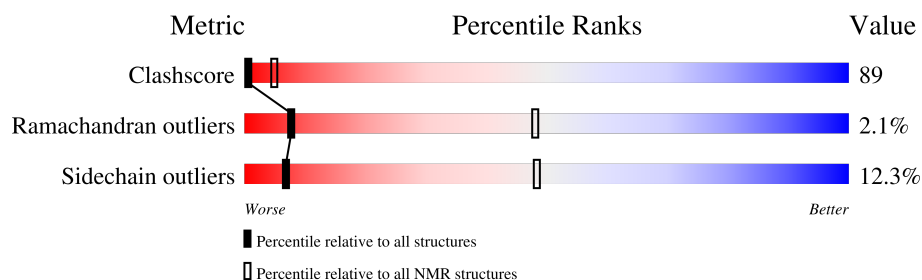
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

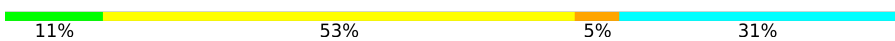
The overall completeness of chemical shifts assignment is 60%.

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	55	

2 Ensemble composition and analysis

This entry contains 21 models. Model 2 is the overall representative, medoid model (most similar to other models). The authors have identified model 3 as representative.

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:11-A:48 (38)	0.48	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 4 clusters and 4 single-model clusters were found.

Cluster number	Models
1	1, 2, 4, 6, 8, 9, 11, 12, 20
2	3, 5, 7
3	10, 13, 17
4	14, 15
Single-model clusters	16; 18; 19; 21

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 892 atoms, of which 443 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called SERINE PROTEINASE INHIBITOR LEKTI.

Mol	Chain	Residues	Atoms						Trace
1	A	55	Total	C	H	N	O	S	0
			892	284	443	74	83	8	

There are 4 discrepancies between the modelled and reference sequences:

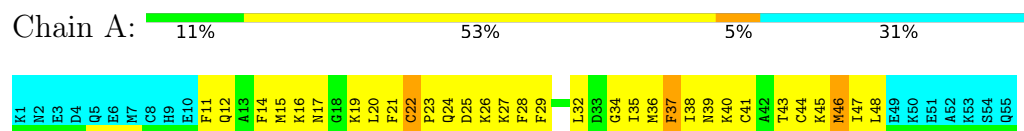
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ASN	ASP	variant	UNP Q9NQ38
A	3	GLU	SER	variant	UNP Q9NQ38
A	4	ASP	LEU	variant	UNP Q9NQ38
A	5	GLN	SER	variant	UNP Q9NQ38

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.

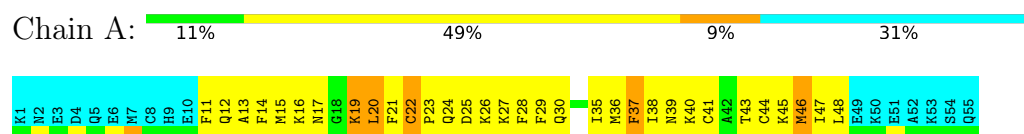
- Molecule 1: SERINE PROTEINASE INHIBITOR LEKTI



4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 2. Colouring as in section 4.1 above.

- Molecule 1: SERINE PROTEINASE INHIBITOR LEKTI



5 Refinement protocol and experimental data overview

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

Of the 60 calculated structures, 21 were deposited, based on the following criterion: *LOWEST ENERGY*.

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

Software name	Classification	Version
X-PLOR	refinement	3.851
NDEE	structure solution	
NMRView	structure solution	4.1.0
X-PLOR	structure solution	3.851

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	2
Total number of shifts	737
Number of shifts mapped to atoms	737
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	60%

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	1.37±0.01	0±0/314 (0.0± 0.0%)	1.57±0.01	0±0/416 (0.0± 0.1%)
All	All	1.37	0/6594 (0.0%)	1.57	1/8736 (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	25	ASP	N-CA-CB	-5.05	104.70	110.35	12	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	307	316	316	55±6
All	All	6447	6636	6636	1163

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 89.

5 of 229 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:20:LEU:HD13	1:A:20:LEU:O	0.83	1.72	18	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:29:PHE:CD2	1:A:35:ILE:HD11	0.82	2.10	5	3
1:A:29:PHE:CE1	1:A:35:ILE:HD11	0.79	2.11	6	10
1:A:29:PHE:CZ	1:A:35:ILE:HD11	0.78	2.12	6	12
1:A:20:LEU:HD23	1:A:45:LYS:HA	0.78	1.55	14	3

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	38/55 (69%)	30±1 (80±3%)	7±1 (18±4%)	1±1 (2±2%)	8	48
All	All	798/1155 (69%)	636 (80%)	145 (18%)	17 (2%)	8	48

All 5 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	23	PRO	8
1	A	41	CYS	4
1	A	25	ASP	2
1	A	27	LYS	2
1	A	29	PHE	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	34/50 (68%)	30±1 (88±3%)	4±1 (12±3%)	7	48
All	All	714/1050 (68%)	626 (88%)	88 (12%)	7	48

5 of 12 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	37	PHE	21
1	A	46	MET	20
1	A	22	CYS	13
1	A	35	ILE	9
1	A	20	LEU	8

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 [i](#)

The completeness of assignment taking into account all chemical shift lists is 60% for the well-defined parts and 56% 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 [i](#)

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	367
Number of shifts mapped to atoms	367
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	0	—	None (insufficient data)
$^{13}\text{C}_\beta$	0	—	None (insufficient data)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	46	-0.32 ± 0.64	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 49%, i.e. 266 atoms were assigned a chemical shift out of a possible 541. 0 out of 3 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	104/190 (55%)	70/77 (91%)	0/76 (0%)	34/37 (92%)
Sidechain	146/291 (50%)	142/189 (75%)	0/91 (0%)	4/11 (36%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	16/60 (27%)	16/30 (53%)	0/30 (0%)	0/0 (—%)
Overall	266/541 (49%)	228/296 (77%)	0/197 (0%)	38/48 (79%)

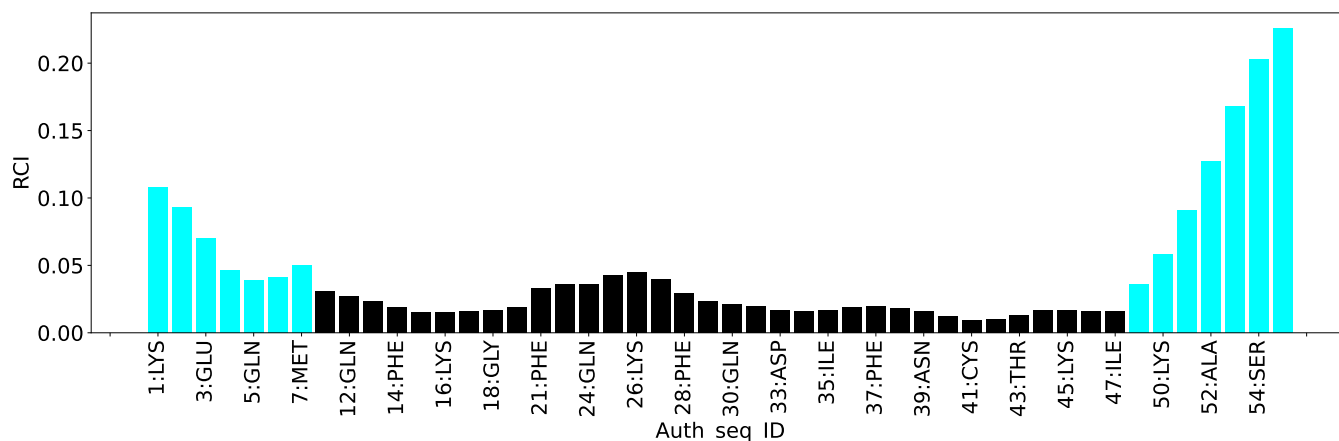
7.1.4 Statistically unusual chemical shifts [i](#)

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:



7.2 Chemical shift list 2

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_2*

7.2.1 Bookkeeping [i](#)

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	370
Number of shifts mapped to atoms	370
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

7.2.2 Chemical shift referencing [i](#)

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

7.2.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 52%, i.e. 281 atoms were assigned a chemical shift out of a possible 541. 0 out of 3 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	77/190 (41%)	77/77 (100%)	0/76 (0%)	0/37 (0%)
Sidechain	178/291 (61%)	178/189 (94%)	0/91 (0%)	0/11 (0%)
Aromatic	26/60 (43%)	26/30 (87%)	0/30 (0%)	0/0 (—%)
Overall	281/541 (52%)	281/296 (95%)	0/197 (0%)	0/48 (0%)

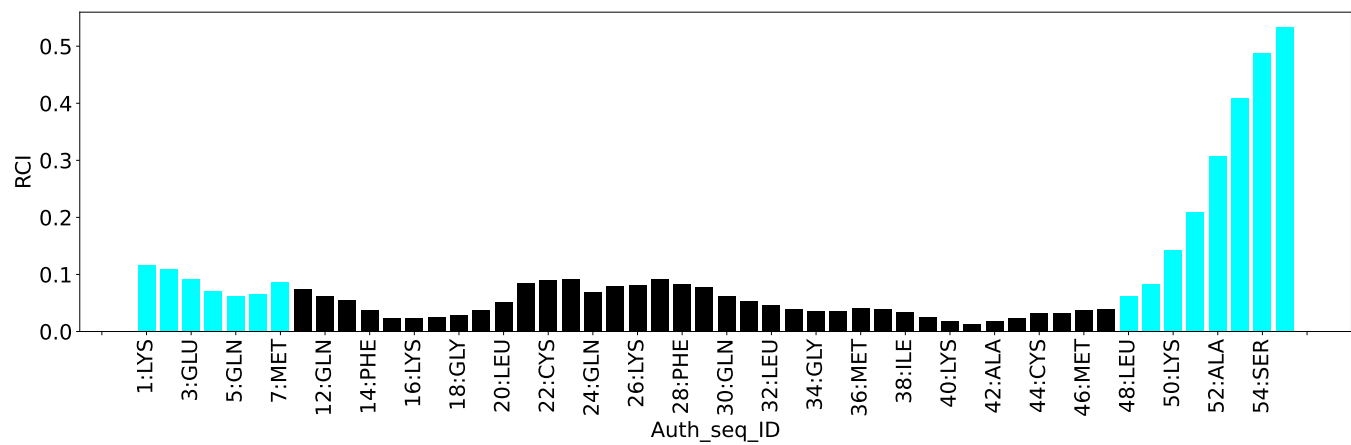
7.2.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.2.5 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 [i](#)

8.1 Conformationally restricting restraints [i](#)

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	523
Intra-residue ($ i-j =0$)	102
Sequential ($ i-j =1$)	203
Medium range ($ i-j >1$ and $ i-j <5$)	152
Long range ($ i-j \geq 5$)	66
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	14
Number of unmapped restraints	0
Number of restraints per residue	9.8
Number of long range restraints per residue ¹	1.2

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 [i](#)

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 [i](#)

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)	2.0	0.2
0.2-0.5 (Medium)	4.1	0.5
>0.5 (Large)	15.7	3.38

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

Dihedral-angle violations less than 1° are not included in the calculation. There are no dihedral-angle violations

9 Distance violation analysis [i](#)

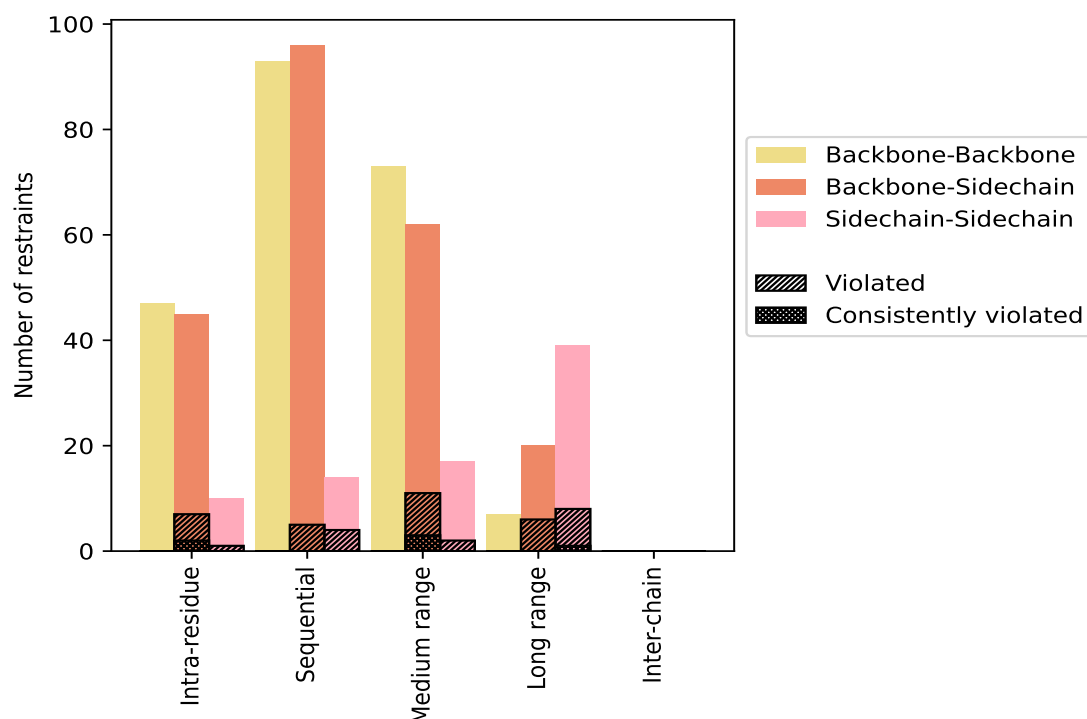
9.1 Summary of distance violations [i](#)

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	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	102	19.5	8	7.8	1.5	2	2.0	0.4
Backbone-Backbone	47	9.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	45	8.6	7	15.6	1.3	2	4.4	0.4
Sidechain-Sidechain	10	1.9	1	10.0	0.2	0	0.0	0.0
Sequential ($i-j =1$)	203	38.8	9	4.4	1.7	0	0.0	0.0
Backbone-Backbone	93	17.8	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	96	18.4	5	5.2	1.0	0	0.0	0.0
Sidechain-Sidechain	14	2.7	4	28.6	0.8	0	0.0	0.0
Medium range ($i-j >1$ & $i-j <5$)	152	29.1	13	8.6	2.5	3	2.0	0.6
Backbone-Backbone	73	14.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	62	11.9	11	17.7	2.1	3	4.8	0.6
Sidechain-Sidechain	17	3.3	2	11.8	0.4	0	0.0	0.0
Long range ($i-j \geq 5$)	66	12.6	14	21.2	2.7	1	1.5	0.2
Backbone-Backbone	7	1.3	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	20	3.8	6	30.0	1.1	0	0.0	0.0
Sidechain-Sidechain	39	7.5	8	20.5	1.5	1	2.6	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	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	523	100.0	44	8.4	8.4	6	1.1	1.1
Backbone-Backbone	220	42.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	223	42.6	29	13.0	5.5	5	2.2	1.0
Sidechain-Sidechain	80	15.3	15	18.8	2.9	1	1.2	0.2

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	4	4	7	11	0	26	0.88	2.73	0.69	0.72
2	6	3	6	13	0	28	0.97	3.1	0.72	0.74
3	3	3	6	13	0	25	0.93	3.38	0.77	0.72
4	6	4	8	10	0	28	0.98	2.81	0.69	0.77
5	4	4	8	11	0	27	0.92	2.98	0.69	0.73
6	6	2	9	4	0	21	0.97	1.96	0.51	0.79
7	4	4	5	8	0	21	0.77	1.94	0.42	0.71
8	6	3	8	11	0	28	1.02	3.1	0.76	0.72
9	4	4	6	13	0	27	1.0	2.49	0.64	0.87
10	2	0	5	1	0	8	0.7	1.51	0.39	0.59

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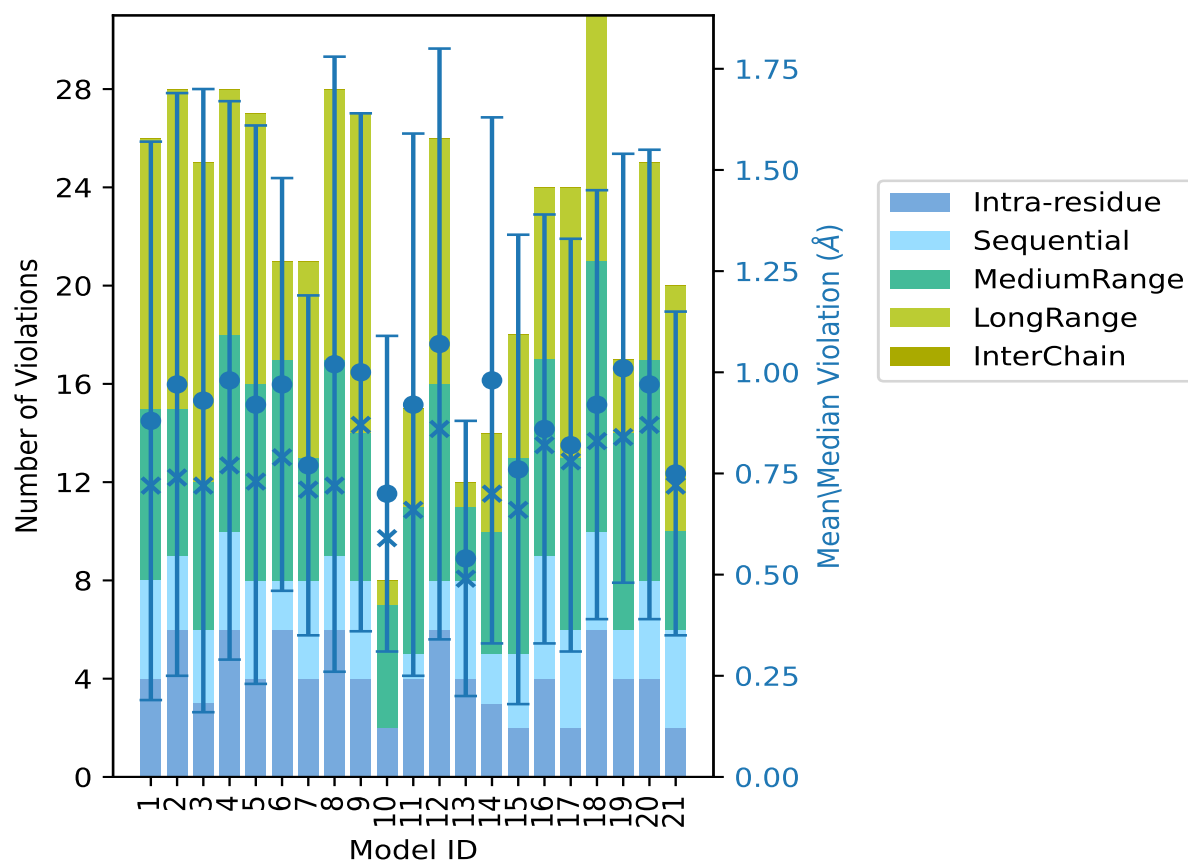
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
11	4	1	6	4	0	15	0.92	2.03	0.67	0.66
12	6	2	8	10	0	26	1.07	3.17	0.73	0.86
13	4	4	3	1	0	12	0.54	1.3	0.34	0.49
14	3	2	5	4	0	14	0.98	2.03	0.65	0.7
15	2	3	8	5	0	18	0.76	1.9	0.58	0.66
16	4	5	8	7	0	24	0.86	2.07	0.53	0.82
17	2	4	7	11	0	24	0.82	2.29	0.51	0.78
18	6	4	11	10	0	31	0.92	2.17	0.53	0.83
19	4	2	8	3	0	17	1.01	1.95	0.53	0.84
20	4	4	9	8	0	25	0.97	2.31	0.58	0.87
21	2	4	4	10	0	20	0.75	1.4	0.4	0.72

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 ⓘ



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](#)

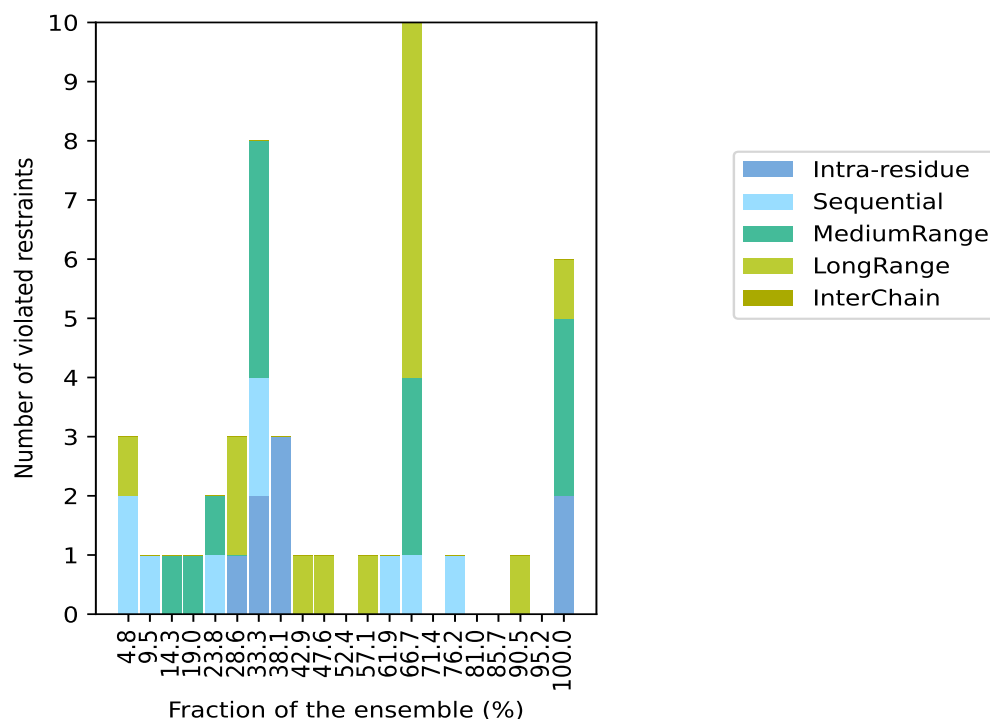
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 479(IR:94, SQ:194, MR:139, LR:52, IC:0) restraints are not 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	2	0	1	0	3	1	4.8
0	1	0	0	0	1	2	9.5
0	0	1	0	0	1	3	14.3
0	0	1	0	0	1	4	19.0
0	1	1	0	0	2	5	23.8
1	0	0	2	0	3	6	28.6
2	2	4	0	0	8	7	33.3
3	0	0	0	0	3	8	38.1
0	0	0	1	0	1	9	42.9
0	0	0	1	0	1	10	47.6
0	0	0	0	0	0	11	52.4
0	0	0	1	0	1	12	57.1
0	1	0	0	0	1	13	61.9
0	1	3	6	0	10	14	66.7
0	0	0	0	0	0	15	71.4
0	1	0	0	0	1	16	76.2
0	0	0	0	0	0	17	81.0
0	0	0	0	0	0	18	85.7
0	0	0	1	0	1	19	90.5
0	0	0	0	0	0	20	95.2
2	0	3	1	0	6	21	100.0

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

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

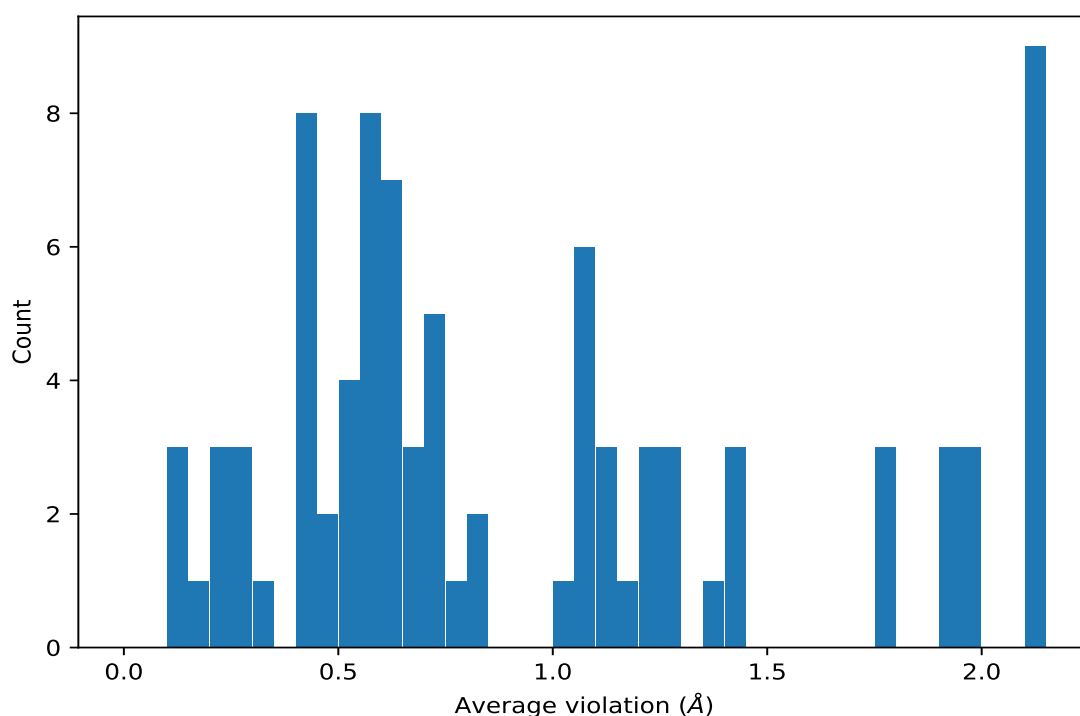
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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)
(1,451)	1:34:A:GLY:HA3	1:37:A:PHE:HB2	21	1.38	0.09	1.39
(1,316)	1:40:A:LYS:H	1:36:A:MET:HB2	21	1.07	0.07	1.08
(1,433)	1:15:A:MET:HB2	1:20:A:LEU:HB3	21	1.04	0.71	0.79
(1,70)	1:36:A:MET:H	1:36:A:MET:HB3	21	0.72	0.06	0.72
(1,92)	1:47:A:ILE:H	1:47:A:ILE:HG13	21	0.71	0.04	0.72
(1,314)	1:39:A:ASN:H	1:36:A:MET:HB2	21	0.4	0.07	0.39
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	19	2.1	0.81	1.83
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	19	2.1	0.81	1.83
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	19	2.1	0.81	1.83
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	19	2.1	0.81	1.83
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	19	2.1	0.81	1.83
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	19	2.1	0.81	1.83
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	19	2.1	0.81	1.83
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	19	2.1	0.81	1.83
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	19	2.1	0.81	1.83
(1,415)	1:49:A:GLU:HG3	1:48:A:LEU:HD11	16	1.14	0.62	0.74

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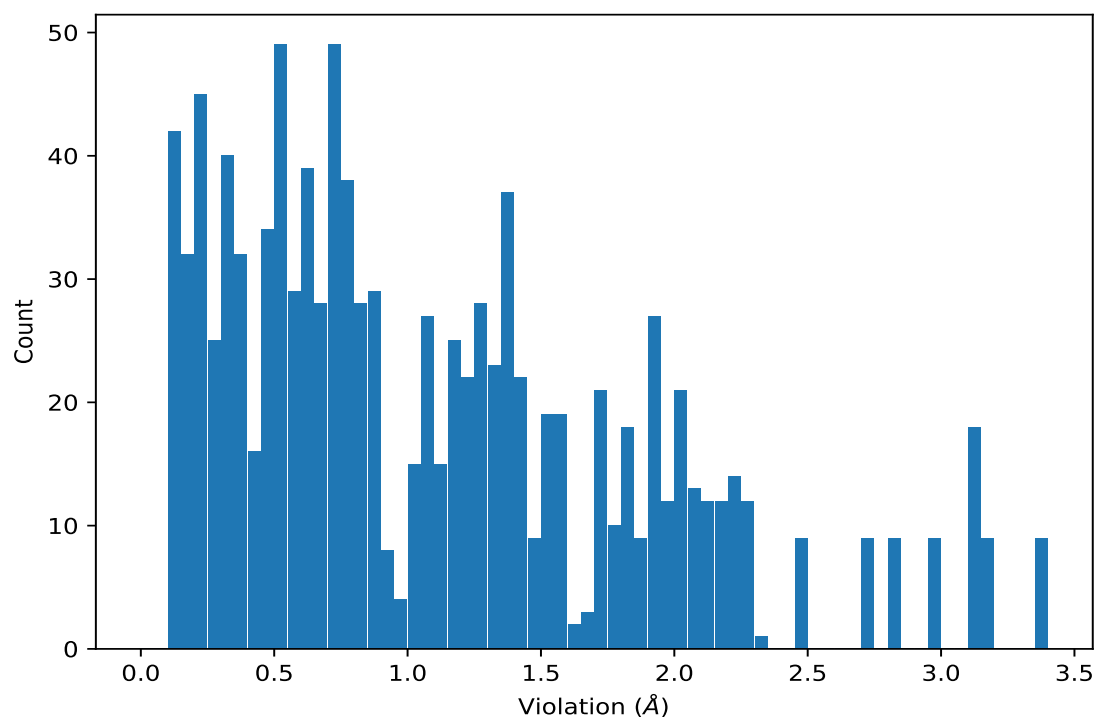
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,415)	1:49:A:GLU:HG3	1:48:A:LEU:HD12	16	1.14	0.62	0.74
(1,415)	1:49:A:GLU:HG3	1:48:A:LEU:HD13	16	1.14	0.62	0.74
(1,475)	1:46:A:MET:HG3	1:48:A:LEU:HD11	14	1.97	0.21	1.98
(1,475)	1:46:A:MET:HG3	1:48:A:LEU:HD12	14	1.97	0.21	1.98
(1,475)	1:46:A:MET:HG3	1:48:A:LEU:HD13	14	1.97	0.21	1.98
(1,147)	1:21:A:PHE:H	1:20:A:LEU:HD11	14	1.91	0.56	2.13

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)
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	3	3.38
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	3	3.38
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	3	3.38
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	3	3.38
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	3	3.38
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	3	3.38
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	3	3.38
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	3	3.38
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	3	3.38
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	12	3.17
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	12	3.17
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	12	3.17
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	12	3.17
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	12	3.17
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	12	3.17
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	12	3.17
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	12	3.17
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	12	3.17
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	2	3.1
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	2	3.1
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	2	3.1
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	2	3.1
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	2	3.1
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	2	3.1
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	2	3.1
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	2	3.1
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	2	3.1
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	8	3.1
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	8	3.1
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	8	3.1
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	8	3.1
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	8	3.1
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	8	3.1
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	8	3.1
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	8	3.1
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	8	3.1
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	5	2.98
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	5	2.98
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	5	2.98
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	5	2.98
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	5	2.98
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	5	2.98
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	5	2.98

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	5	2.98
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	5	2.98
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	4	2.81
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	4	2.81
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	4	2.81
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	4	2.81
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	4	2.81
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	4	2.81
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	4	2.81
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	4	2.81
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	4	2.81
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	1	2.73
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	1	2.73
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	1	2.73
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	1	2.73
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	1	2.73
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	1	2.73
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	1	2.73
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	1	2.73
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	1	2.73
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD21	9	2.49
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD22	9	2.49
(1,523)	1:48:A:LEU:HD11	1:20:A:LEU:HD23	9	2.49
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD21	9	2.49
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD22	9	2.49
(1,523)	1:48:A:LEU:HD12	1:20:A:LEU:HD23	9	2.49
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD21	9	2.49
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD22	9	2.49
(1,523)	1:48:A:LEU:HD13	1:20:A:LEU:HD23	9	2.49
(1,433)	1:15:A:MET:HB2	1:20:A:LEU:HB3	20	2.31
(1,147)	1:21:A:PHE:H	1:20:A:LEU:HD11	17	2.29

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

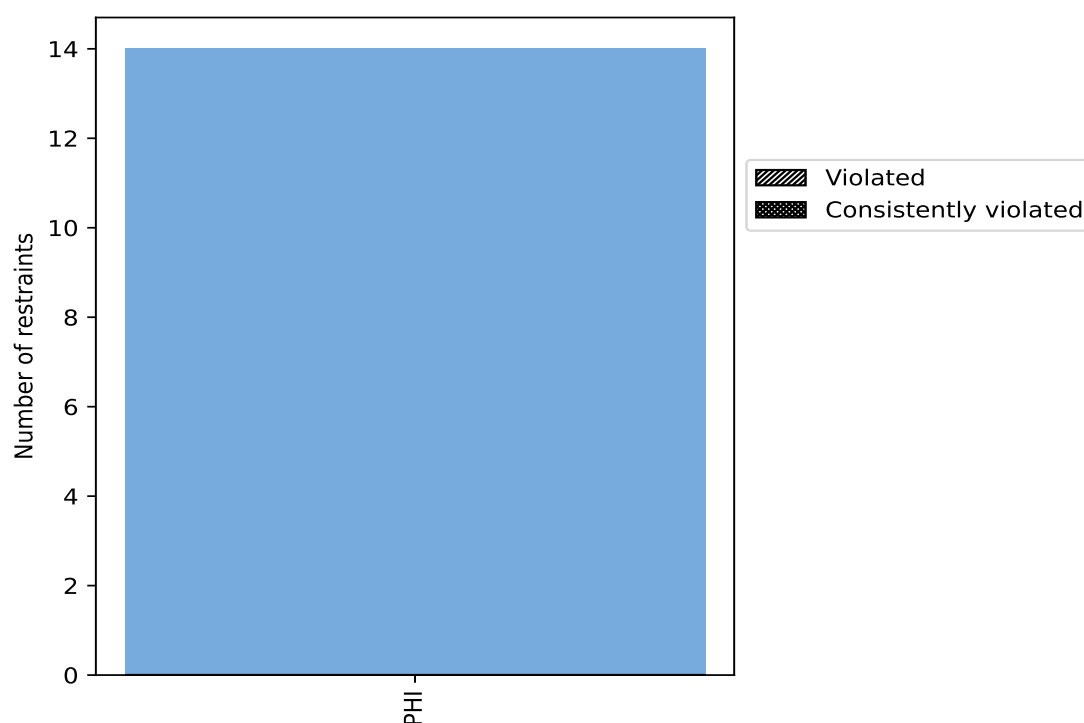
10.1 Summary of dihedral-angle violations [i](#)

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	14	100.0	0	0.0	0.0	0	0.0	0.0
Total	14	100.0	0	0.0	0.0	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 [i](#)

No violations found

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

No violations found

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

No violations found

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

No violations found