



Full wwPDB NMR Structure Validation Report ⓘ

Mar 26, 2026 – 07:51 AM UTC

PDB ID : 2M8G / pdb_00002m8g
BMRB ID : 19251
Title : Structure, function, and tethering of DNA-binding domains in 54 transcriptional activators
Authors : Hong, E.; Wemmer, D.
Deposited on : 2013-05-19

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 Phenix2.0
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

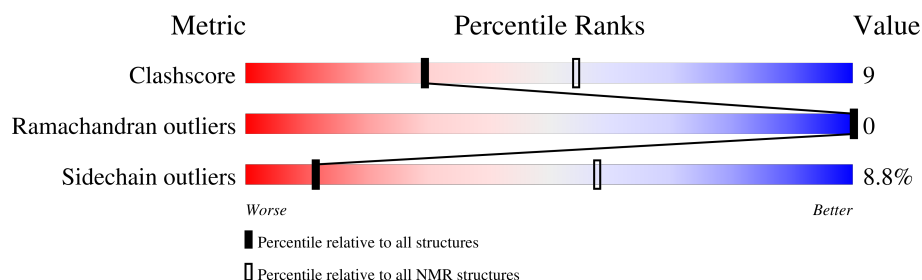
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 75%.

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	X	70	43% 14% 43%

2 Ensemble composition and analysis

This entry contains 20 models. Model 11 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *fewest violations*.

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	X:454-X:493 (40)	0.11	11

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 2 clusters and 5 single-model clusters were found.

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

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1101 atoms, of which 515 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Transcriptional regulator.

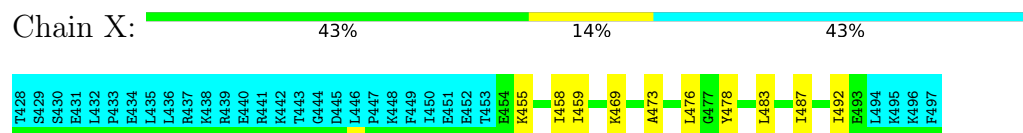
Mol	Chain	Residues	Atoms					Trace
1	X	70	Total	C	H	N	O	0
			1101	377	515	104	105	

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: Transcriptional regulator

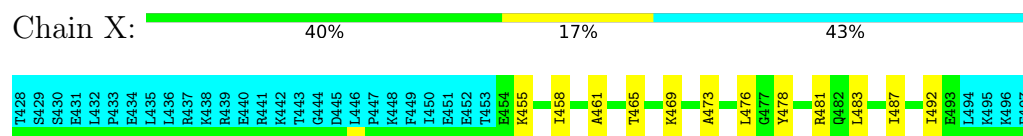


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

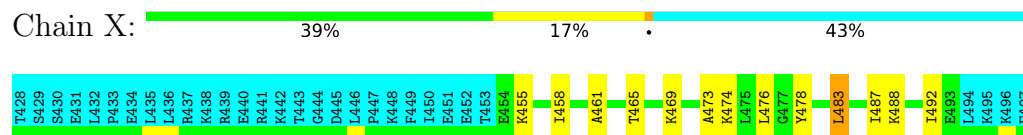
4.2.1 Score per residue for model 1

- Molecule 1: Transcriptional regulator



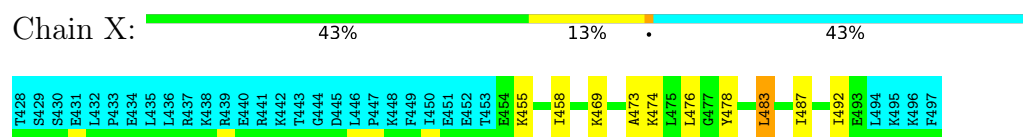
4.2.2 Score per residue for model 2

- Molecule 1: Transcriptional regulator



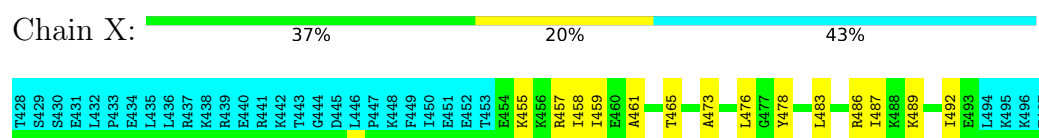
4.2.3 Score per residue for model 3

- Molecule 1: Transcriptional regulator



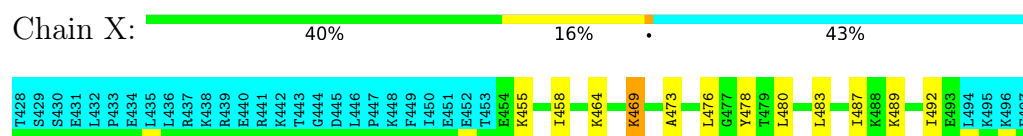
4.2.4 Score per residue for model 4

- Molecule 1: Transcriptional regulator



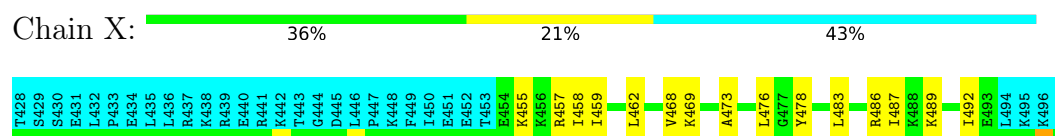
4.2.5 Score per residue for model 5

- Molecule 1: Transcriptional regulator



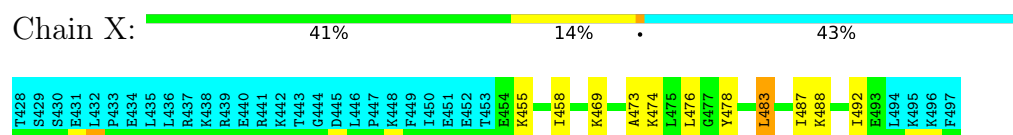
4.2.6 Score per residue for model 6

- Molecule 1: Transcriptional regulator



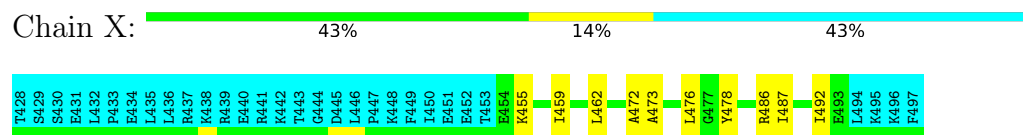
4.2.7 Score per residue for model 7

- Molecule 1: Transcriptional regulator



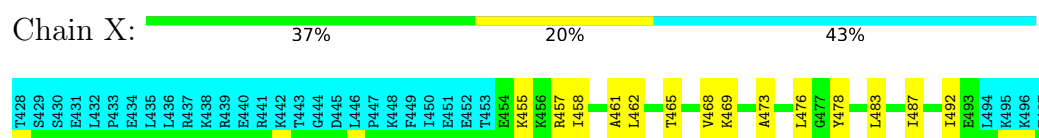
4.2.8 Score per residue for model 8

- Molecule 1: Transcriptional regulator



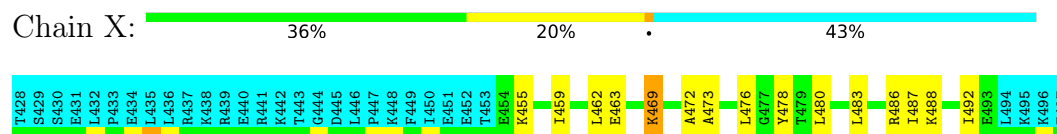
4.2.9 Score per residue for model 9

- Molecule 1: Transcriptional regulator



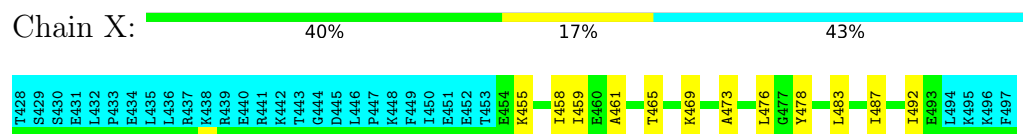
4.2.10 Score per residue for model 10

- Molecule 1: Transcriptional regulator



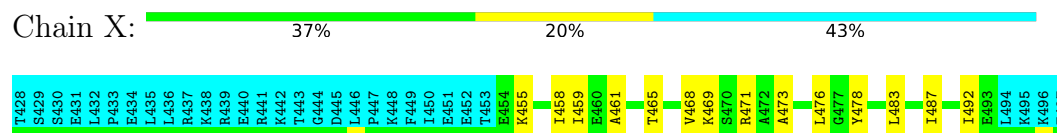
4.2.11 Score per residue for model 11 (medoid)

- Molecule 1: Transcriptional regulator



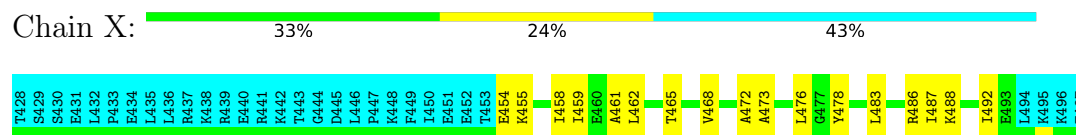
4.2.12 Score per residue for model 12

- Molecule 1: Transcriptional regulator



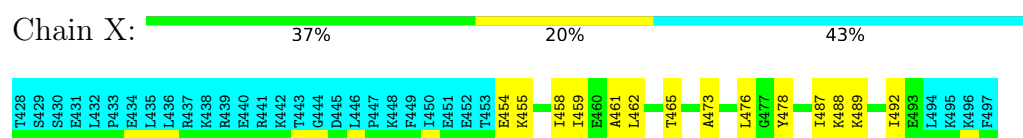
4.2.13 Score per residue for model 13

- Molecule 1: Transcriptional regulator



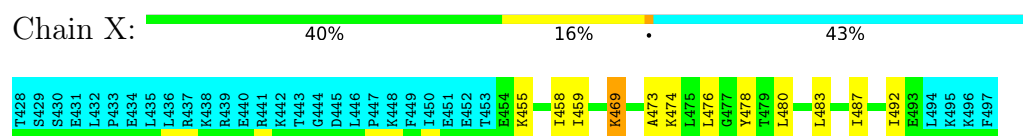
4.2.14 Score per residue for model 14

- Molecule 1: Transcriptional regulator



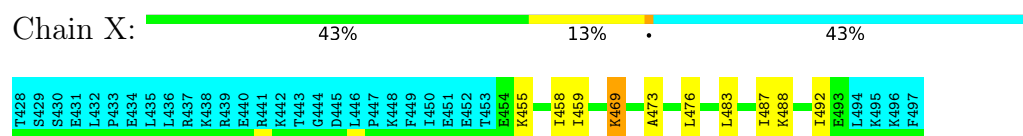
4.2.15 Score per residue for model 15

- Molecule 1: Transcriptional regulator



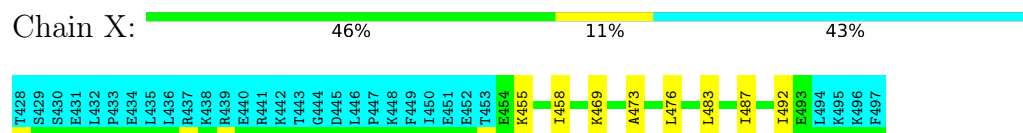
4.2.16 Score per residue for model 16

- Molecule 1: Transcriptional regulator



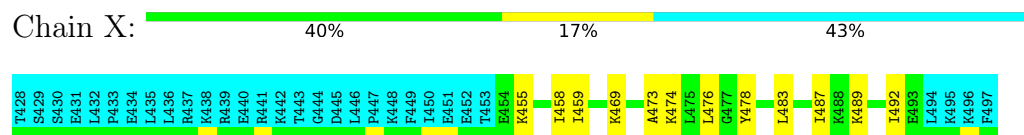
4.2.17 Score per residue for model 17

- Molecule 1: Transcriptional regulator



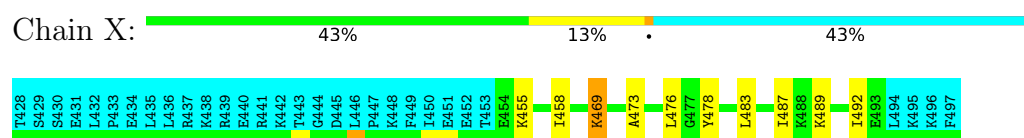
4.2.18 Score per residue for model 18

- Molecule 1: Transcriptional regulator



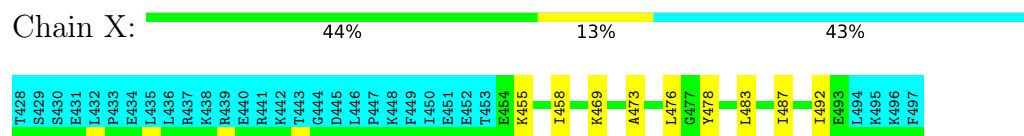
4.2.19 Score per residue for model 19

- Molecule 1: Transcriptional regulator



4.2.20 Score per residue for model 20

- Molecule 1: Transcriptional regulator



5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *target function*.

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

Software name	Classification	Version
CYANA	structure solution	
MOLMOL	refinement	
CYANA	refinement	

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

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	658
Number of shifts mapped to atoms	577
Number of unparsed shifts	0
Number of shifts with mapping errors	81
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	75%

6 Model quality

6.1 Standard geometry

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	X	334	301	366	6±1
All	All	6680	6020	7320	129

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 9.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:X:458:ILE:HD13	1:X:476:LEU:HD13	0.67	1.65	16	18
1:X:469:LYS:HG3	1:X:483:LEU:HD23	0.65	1.67	6	1
1:X:487:ILE:HD13	1:X:492:ILE:HD12	0.64	1.69	12	20
1:X:459:ILE:HD12	1:X:492:ILE:HG23	0.59	1.73	10	11
1:X:469:LYS:CE	1:X:480:LEU:HD13	0.58	2.28	15	1
1:X:473:ALA:HB2	1:X:483:LEU:HD22	0.56	1.78	13	3
1:X:469:LYS:CG	1:X:483:LEU:HD23	0.55	2.31	6	2
1:X:473:ALA:HA	1:X:483:LEU:HD13	0.55	1.79	12	13
1:X:473:ALA:HB1	1:X:478:TYR:O	0.55	2.02	13	18
1:X:469:LYS:HG2	1:X:483:LEU:HD23	0.51	1.82	15	2
1:X:480:LEU:HD12	1:X:480:LEU:O	0.51	2.06	15	1
1:X:469:LYS:O	1:X:483:LEU:HD22	0.51	2.06	16	13
1:X:472:ALA:O	1:X:476:LEU:HD12	0.51	2.06	10	3
1:X:469:LYS:HE2	1:X:480:LEU:HD13	0.49	1.83	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:X:473:ALA:CA	1:X:483:LEU:HD13	0.47	2.38	2	7
1:X:461:ALA:O	1:X:465:THR:HG23	0.47	2.09	1	8
1:X:462:LEU:HD23	1:X:468:VAL:O	0.47	2.10	9	3
1:X:468:VAL:HG21	1:X:471:ARG:CZ	0.46	2.40	12	1
1:X:462:LEU:HD11	1:X:487:ILE:HD11	0.46	1.88	8	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	X	40/70 (57%)	40±0 (100±0%)	0±0 (0±0%)	0±0 (0±0%)	100	100
All	All	800/1400 (57%)	800 (100%)	0 (0%)	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	X	34/63 (54%)	31±1 (91±4%)	3±1 (9±4%)	11	58
All	All	680/1260 (54%)	620 (91%)	60 (9%)	11	58

All 13 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	X	455	LYS	20
1	X	488	LYS	6
1	X	489	LYS	6

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Mol	Chain	Res	Type	Models (Total)
1	X	474	LYS	5
1	X	486	ARG	5
1	X	469	LYS	5
1	X	483	LEU	3
1	X	457	ARG	3
1	X	480	LEU	2
1	X	454	GLU	2
1	X	481	ARG	1
1	X	464	LYS	1
1	X	463	GLU	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no 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 75% for the well-defined parts and 62% 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	658
Number of shifts mapped to atoms	577
Number of unparsed shifts	0
Number of shifts with mapping errors	81
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. All 81 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	X	430	SER	HB2	3.898	0.02	1
1	X	431	GLU	HB2	1.938	0.02	1
1	X	431	GLU	HG2	2.254	0.02	2
1	X	432	LEU	HB2	1.619	0.02	1
1	X	433	PRO	HB2	2.33	0.02	2
1	X	433	PRO	HD2	3.897	0.02	2
1	X	433	PRO	HG2	2.062	0.02	1
1	X	434	GLU	HB2	2.031	0.02	1
1	X	445	ASP	HB2	2.722	0.02	2
1	X	448	LYS	HB2	1.789	0.02	1
1	X	449	PHE	HB2	3.198	0.02	1
1	X	450	ILE	HG12	1.501	0.02	2
1	X	451	GLU	HB2	2.113	0.02	1
1	X	451	GLU	HG2	2.393	0.02	2

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	X	452	GLU	HB2	2.093	0.02	2
1	X	454	GLU	HB2	2.123	0.02	1
1	X	454	GLU	HG2	2.478	0.02	1
1	X	455	LYS	HB2	1.68	0.02	2
1	X	455	LYS	HD2	1.683	0.02	1
1	X	455	LYS	HE2	2.957	0.02	2
1	X	455	LYS	HG2	1.204	0.02	2
1	X	456	LYS	HB2	1.897	0.02	1
1	X	456	LYS	HG2	1.573	0.02	2
1	X	457	ARG	HB2	1.978	0.02	2
1	X	457	ARG	HD2	3.234	0.02	2
1	X	457	ARG	HG2	1.77	0.02	2
1	X	458	ILE	HG12	2.144	0.02	2
1	X	459	ILE	HG12	1.847	0.02	2
1	X	460	GLU	HB2	2.122	0.02	2
1	X	460	GLU	HG2	2.465	0.02	2
1	X	462	LEU	HB2	1.693	0.02	2
1	X	463	GLU	HB2	2.235	0.02	2
1	X	463	GLU	HG2	2.542	0.02	2
1	X	464	LYS	HB2	1.957	0.02	2
1	X	464	LYS	HD2	1.668	0.02	1
1	X	464	LYS	HG2	1.616	0.02	2
1	X	467	TYR	HB2	3.343	0.02	2
1	X	471	ARG	HB2	1.993	0.02	2
1	X	471	ARG	HD2	3.281	0.02	1
1	X	471	ARG	HG2	1.833	0.02	2
1	X	474	LYS	HB2	1.984	0.02	1
1	X	474	LYS	HD2	1.755	0.02	1
1	X	474	LYS	HG2	1.623	0.02	2
1	X	475	LEU	HB2	1.814	0.02	2
1	X	476	LEU	HB2	1.378	0.02	2
1	X	478	TYR	HB2	3.7	0.02	2
1	X	480	LEU	HB2	1.721	0.02	1
1	X	481	ARG	HB2	1.874	0.02	2
1	X	481	ARG	HD2	3.214	0.02	1
1	X	481	ARG	HG2	1.735	0.02	2
1	X	482	GLN	HB2	2.326	0.02	2
1	X	482	GLN	HG2	2.364	0.02	2
1	X	483	LEU	HB2	2.26	0.02	2
1	X	484	ASP	HB2	2.767	0.02	2
1	X	485	TYR	HB2	3.245	0.02	2

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	X	486	ARG	HB2	1.764	0.02	1
1	X	486	ARG	HD2	3.04	0.02	2
1	X	486	ARG	HG2	1.904	0.02	2
1	X	487	ILE	HG12	1.891	0.02	2
1	X	488	LYS	HB2	1.852	0.02	2
1	X	488	LYS	HD2	1.674	0.02	1
1	X	488	LYS	HE2	2.968	0.02	1
1	X	488	LYS	HG2	1.468	0.02	2
1	X	489	LYS	HB2	1.259	0.02	2
1	X	489	LYS	HD2	1.393	0.02	2
1	X	489	LYS	HE2	2.843	0.02	1
1	X	489	LYS	HG2	1.086	0.02	2
1	X	490	TYR	HB2	3.265	0.02	2
1	X	492	ILE	HG12	1.703	0.02	2
1	X	493	GLU	HB2	1.942	0.02	1
1	X	493	GLU	HG2	2.267	0.02	2
1	X	494	LEU	HB2	1.739	0.02	2
1	X	495	LYS	HB2	1.258	0.02	2
1	X	495	LYS	HD2	1.48	0.02	1
1	X	495	LYS	HE2	2.836	0.02	1
1	X	495	LYS	HG2	0.9721	0.02	1
1	X	496	LYS	HB2	1.716	0.02	2
1	X	496	LYS	HD2	1.641	0.02	1
1	X	496	LYS	HE2	2.966	0.02	1
1	X	496	LYS	HG2	1.313	0.02	1
1	X	497	PHE	HB2	3.198	0.02	2

7.1.2 Chemical shift referencing ⓘ

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	61	-1.65 ± 0.12	Should be checked
$^{13}\text{C}_\beta$	55	-0.82 ± 0.19	Should be checked
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	57	-0.82 ± 0.14	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 75%, i.e. 460 atoms were assigned a chemical

shift out of a possible 610. 0 out of 6 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	157/203 (77%)	79/83 (95%)	39/80 (49%)	39/40 (98%)
Sidechain	287/371 (77%)	200/239 (84%)	86/112 (77%)	1/20 (5%)
Aromatic	16/36 (44%)	16/16 (100%)	0/20 (0%)	0/0 (—%)
Overall	460/610 (75%)	295/338 (87%)	125/212 (59%)	40/60 (67%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 62%, i.e. 658 atoms were assigned a chemical shift out of a possible 1061. 0 out of 11 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	234/350 (67%)	116/142 (82%)	61/140 (44%)	57/68 (84%)
Sidechain	402/655 (61%)	275/420 (65%)	126/201 (63%)	1/34 (3%)
Aromatic	22/56 (39%)	22/26 (85%)	0/30 (0%)	0/0 (—%)
Overall	658/1061 (62%)	413/588 (70%)	187/371 (50%)	58/102 (57%)

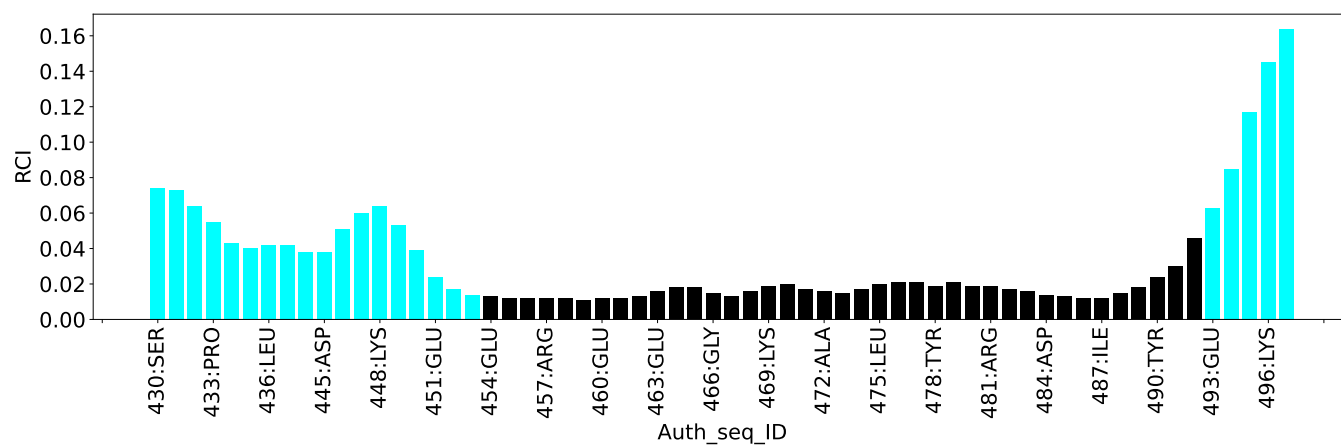
7.1.4 Statistically unusual chemical shifts [i](#)

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 X:



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	1168
Intra-residue ($ i-j =0$)	286
Sequential ($ i-j =1$)	274
Medium range ($ i-j >1$ and $ i-j <5$)	299
Long range ($ i-j \geq 5$)	252
Inter-chain	0
Hydrogen bond restraints	57
Disulfide bond restraints	0
Total dihedral-angle restraints	86
Number of unmapped restraints	0
Number of restraints per residue	17.9
Number of long range restraints per residue ¹	3.6

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)	6.7	0.2
0.2-0.5 (Medium)	9.6	0.35
>0.5 (Large)	None	None

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)	6.1	2.93
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

9 Distance violation analysis ⓘ

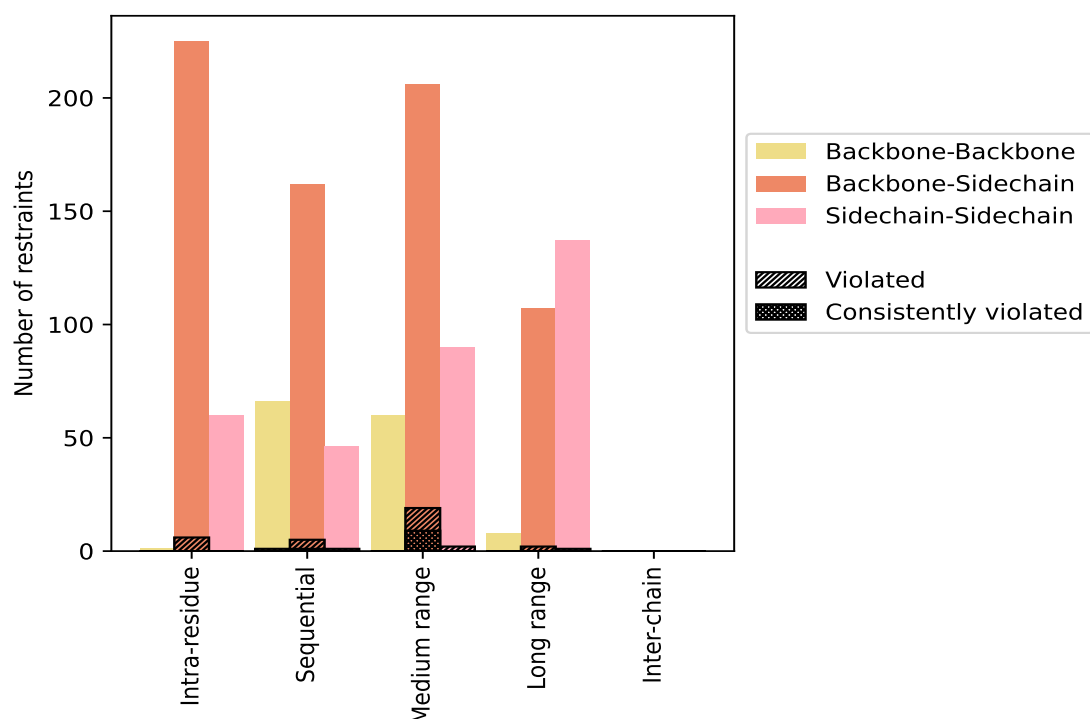
9.1 Summary of distance violations ⓘ

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 ($i-j =0$)	286	24.5	6	2.1	0.5	0	0.0	0.0
Backbone-Backbone	1	0.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	225	19.3	6	2.7	0.5	0	0.0	0.0
Sidechain-Sidechain	60	5.1	0	0.0	0.0	0	0.0	0.0
Sequential ($i-j =1$)	274	23.5	7	2.6	0.6	1	0.4	0.1
Backbone-Backbone	66	5.7	1	1.5	0.1	0	0.0	0.0
Backbone-Sidechain	162	13.9	5	3.1	0.4	1	0.6	0.1
Sidechain-Sidechain	46	3.9	1	2.2	0.1	0	0.0	0.0
Medium range ($i-j >1$ & $i-j <5$)	299	25.6	6	2.0	0.5	0	0.0	0.0
Backbone-Backbone	60	5.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	149	12.8	4	2.7	0.3	0	0.0	0.0
Sidechain-Sidechain	90	7.7	2	2.2	0.2	0	0.0	0.0
Long range ($i-j \geq 5$)	252	21.6	3	1.2	0.3	0	0.0	0.0
Backbone-Backbone	8	0.7	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	107	9.2	2	1.9	0.2	0	0.0	0.0
Sidechain-Sidechain	137	11.7	1	0.7	0.1	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	57	4.9	15	26.3	1.3	9	15.8	0.8
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	1168	100.0	37	3.2	3.2	10	0.9	0.9
Backbone-Backbone	135	11.6	1	0.7	0.1	0	0.0	0.0
Backbone-Sidechain	700	59.9	32	4.6	2.7	10	1.4	0.9
Sidechain-Sidechain	333	28.5	4	1.2	0.3	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	1	1	13	1	0	16	0.21	0.31	0.08	0.24
2	1	2	14	1	0	18	0.21	0.32	0.08	0.22
3	1	2	14	1	0	18	0.21	0.32	0.07	0.22
4	1	1	12	0	0	14	0.23	0.32	0.07	0.26
5	2	1	13	1	0	17	0.21	0.32	0.08	0.23
6	4	1	14	2	0	21	0.2	0.34	0.09	0.17
7	1	2	14	1	0	18	0.21	0.32	0.07	0.22
8	1	1	15	2	0	19	0.21	0.35	0.08	0.19
9	1	1	15	1	0	18	0.22	0.33	0.07	0.22
10	1	1	13	1	0	16	0.21	0.32	0.08	0.22

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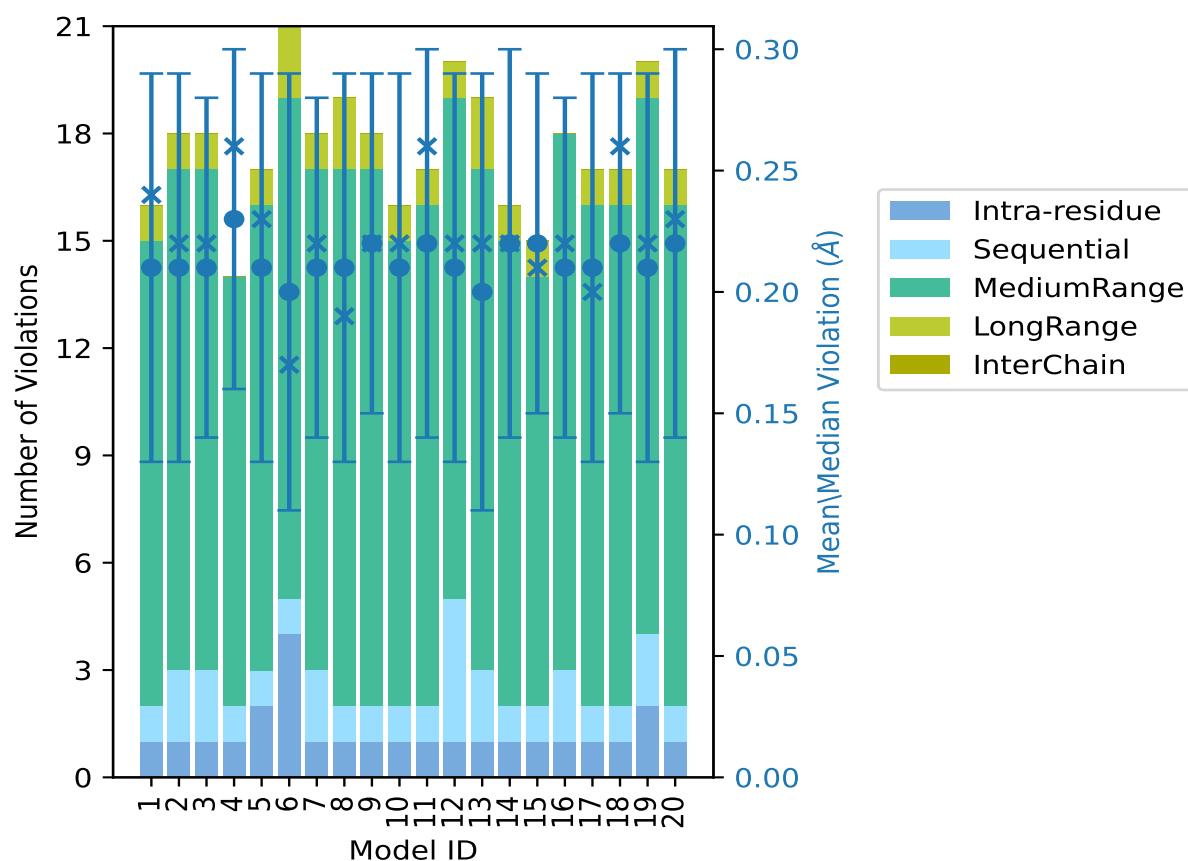
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
11	1	1	14	1	0	17	0.22	0.31	0.08	0.26
12	1	4	14	1	0	20	0.21	0.33	0.08	0.22
13	1	2	14	2	0	19	0.2	0.34	0.09	0.22
14	1	1	13	1	0	16	0.22	0.34	0.08	0.22
15	1	1	12	1	0	15	0.22	0.31	0.07	0.21
16	1	2	15	0	0	18	0.21	0.34	0.07	0.22
17	1	1	14	1	0	17	0.21	0.32	0.08	0.2
18	1	1	14	1	0	17	0.22	0.31	0.07	0.26
19	2	2	15	1	0	20	0.21	0.32	0.08	0.22
20	1	1	14	1	0	17	0.22	0.32	0.08	0.23

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

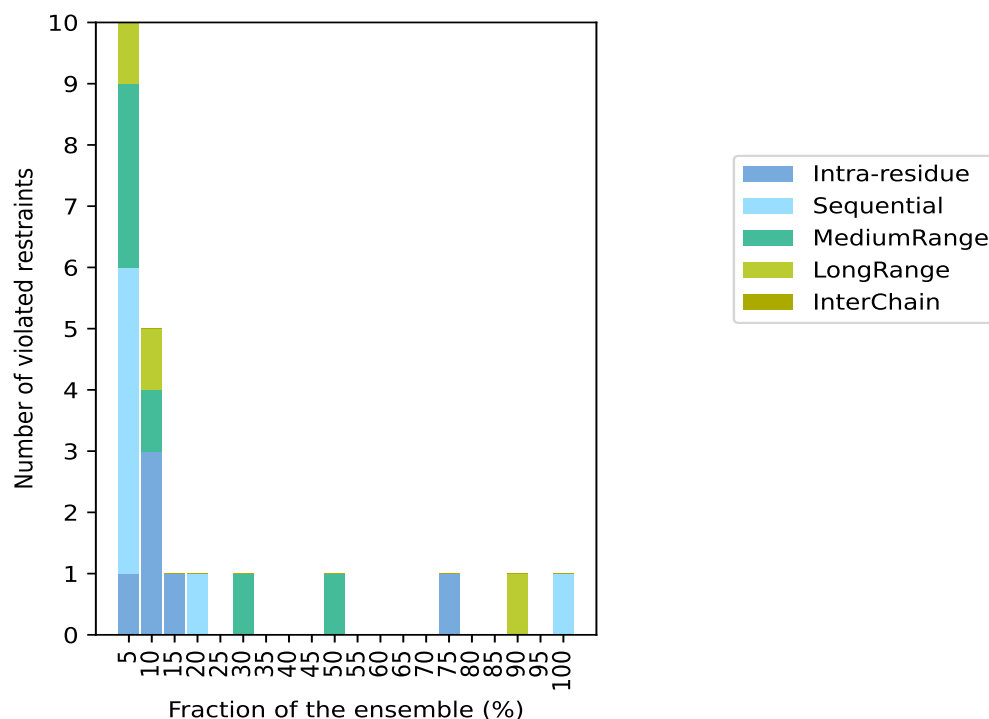
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 1089(IR:280, SQ:267, MR:293, LR:249, IC:0) restraints are 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	%
1	5	3	1	0	10	1	5.0
3	0	1	1	0	5	2	10.0
1	0	0	0	0	1	3	15.0
0	1	0	0	0	1	4	20.0
0	0	0	0	0	0	5	25.0
0	0	1	0	0	1	6	30.0
0	0	0	0	0	0	7	35.0
0	0	0	0	0	0	8	40.0
0	0	0	0	0	0	9	45.0
0	0	1	0	0	1	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
1	0	0	0	0	1	15	75.0
0	0	0	0	0	0	16	80.0
0	0	0	0	0	0	17	85.0
0	0	0	1	0	1	18	90.0
0	0	0	0	0	0	19	95.0
0	1	0	0	0	1	20	100.0

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

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

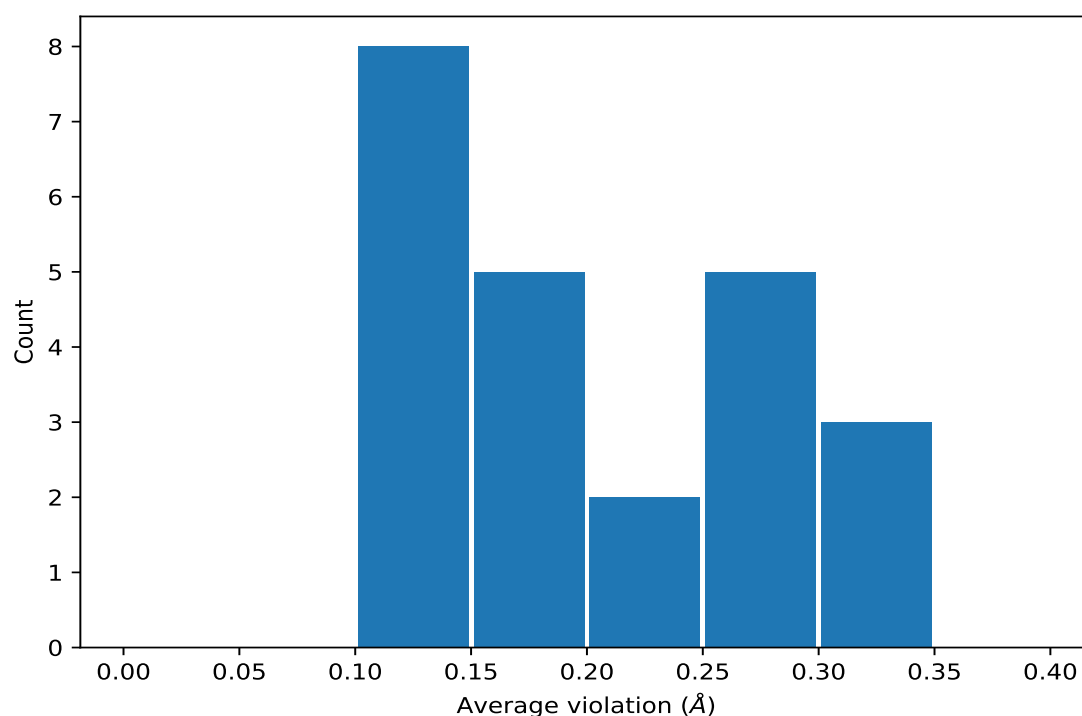
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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)
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	20	0.32	0.01	0.32
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	20	0.31	0.01	0.31
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	20	0.31	0.01	0.31
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	20	0.3	0.02	0.3
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	20	0.29	0.01	0.28
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	20	0.27	0.01	0.27
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	20	0.25	0.02	0.24
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	20	0.23	0.02	0.23
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	20	0.19	0.03	0.2
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	19	0.14	0.01	0.15
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	18	0.12	0.01	0.12
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	17	0.24	0.04	0.25
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	14	0.17	0.04	0.18
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	10	0.14	0.01	0.15
(1,883)	1:432:X:LEU:HB3	1:434:X:GLU:H	6	0.27	0.02	0.26
(2,50)	1:471:X:ARG:O	1:475:X:LEU:N	6	0.17	0.01	0.17

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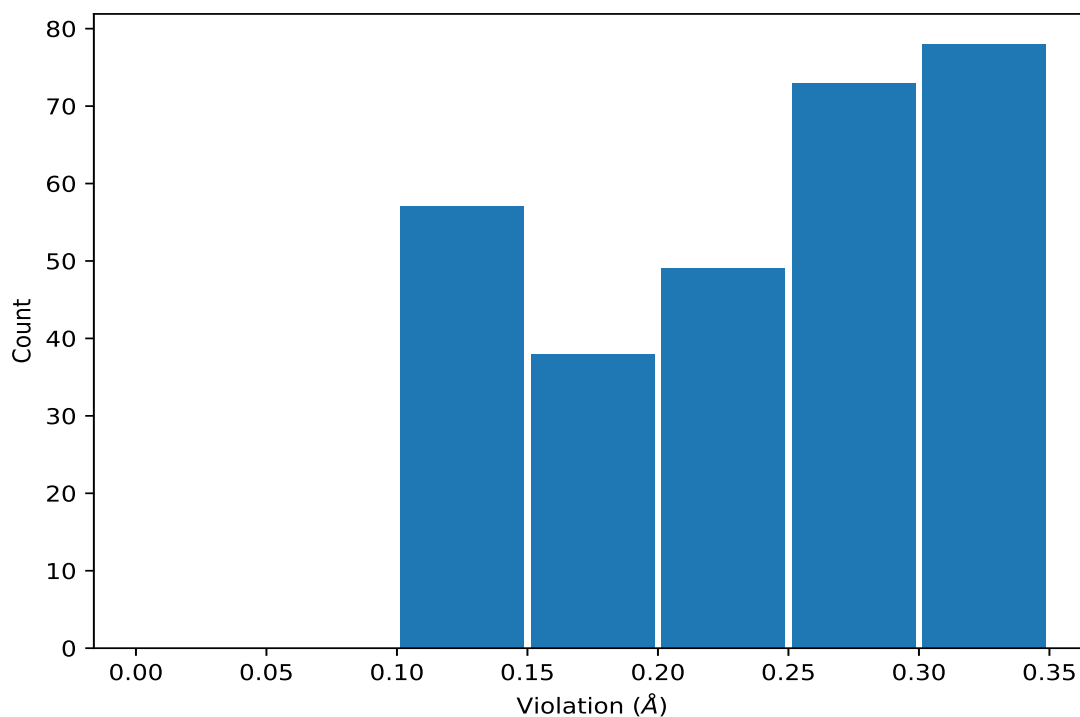
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,1059)	1:486:X:ARG:H	1:486:X:ARG:HG3	3	0.12	0.01	0.13
(1,764)	1:485:X:TYR:H	1:485:X:TYR:HD1	2	0.16	0.02	0.16
(1,764)	1:485:X:TYR:H	1:485:X:TYR:HD2	2	0.16	0.02	0.16
(1,659)	1:472:X:ALA:H	1:474:X:LYS:HD3	2	0.13	0.0	0.13
(1,1071)	1:489:X:LYS:H	1:489:X:LYS:HD3	2	0.12	0.01	0.12
(1,372)	1:458:X:ILE:HA	1:476:X:LEU:HG	2	0.12	0.0	0.12
(1,342)	1:474:X:LYS:HA	1:474:X:LYS:HD3	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)
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	8	0.35
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	6	0.34
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	8	0.34
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	13	0.34
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	14	0.34
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	16	0.34
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	16	0.33
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	6	0.33
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	9	0.33
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	12	0.33
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	4	0.32
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	5	0.32
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	8	0.32
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	14	0.32
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	19	0.32
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	20	0.32
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	2	0.32
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	3	0.32
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	5	0.32
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	6	0.32
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	7	0.32
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	17	0.32
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	19	0.32
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	3	0.32
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	7	0.32
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	17	0.32
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	19	0.32
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	8	0.32
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	10	0.32
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	13	0.32
(1,883)	1:432:X:LEU:HB3	1:434:X:GLU:H	20	0.32
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	1	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	2	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	3	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	7	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	9	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	10	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	11	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	12	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	13	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	16	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	17	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	18	0.31

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	1	0.31
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	9	0.31
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	10	0.31
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	11	0.31
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	12	0.31
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	15	0.31
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	18	0.31
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	20	0.31
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	2	0.31
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	4	0.31
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	10	0.31
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	15	0.31
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	18	0.31
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	2	0.31
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	4	0.31
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	11	0.31
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	14	0.31
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	15	0.31
(1,881)	1:433:X:PRO:HG3	1:434:X:GLU:H	19	0.31
(2,57)	1:484:X:ASP:O	1:488:X:LYS:N	15	0.3
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	4	0.3
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	6	0.3
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	14	0.3
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	1	0.3
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	5	0.3
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	11	0.3
(2,49)	1:470:X:SER:O	1:474:X:LYS:N	20	0.3
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	12	0.3
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	5	0.3
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	12	0.3
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	14	0.3
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	19	0.3
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	6	0.3
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	12	0.3
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	18	0.3
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	13	0.29
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	14	0.29
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	9	0.29
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	10	0.29
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	17	0.29
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	18	0.29
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	1	0.29

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	3	0.29
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	20	0.29
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	7	0.28
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	13	0.28
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	13	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	1	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	3	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	4	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	6	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	7	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	8	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	11	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	16	0.28
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	20	0.28
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	2	0.28
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	5	0.28
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	7	0.28
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	9	0.28
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	16	0.28
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	19	0.28
(2,55)	1:482:X:GLN:O	1:486:X:ARG:N	14	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	1	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	2	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	3	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	8	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	9	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	11	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	12	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	18	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	20	0.27
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	4	0.27
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	5	0.27
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	11	0.27
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	20	0.27
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	11	0.27
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	13	0.27
(2,44)	1:459:X:ILE:O	1:463:X:GLU:N	15	0.27
(2,43)	1:458:X:ILE:O	1:462:X:LEU:N	17	0.27
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	4	0.26
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	5	0.26
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	10	0.26
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	17	0.26

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	13	0.26
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	1	0.26
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	6	0.26
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	18	0.26
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	4	0.26
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	10	0.26
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	18	0.26
(1,883)	1:432:X:LEU:HB3	1:434:X:GLU:H	9	0.26
(1,883)	1:432:X:LEU:HB3	1:434:X:GLU:H	11	0.26
(1,883)	1:432:X:LEU:HB3	1:434:X:GLU:H	15	0.26
(1,883)	1:432:X:LEU:HB3	1:434:X:GLU:H	16	0.26
(1,883)	1:432:X:LEU:HB3	1:434:X:GLU:H	18	0.26
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	15	0.25
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	16	0.25
(2,54)	1:481:X:ARG:O	1:485:X:TYR:N	19	0.25
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	8	0.25
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	2	0.25
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	3	0.25
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	12	0.25
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	1	0.25
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	2	0.25
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	3	0.25
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	9	0.25
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	19	0.25
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	17	0.24
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	7	0.24
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	19	0.24
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	5	0.24
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	13	0.24
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	14	0.24
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	1	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	3	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	4	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	5	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	7	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	9	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	10	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	11	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	12	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	16	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	19	0.23
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	20	0.23

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	6	0.23
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	7	0.23
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	8	0.23
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	16	0.23
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	17	0.23
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	20	0.23
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	8	0.22
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	2	0.22
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	6	0.22
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	18	0.22
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	9	0.22
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	6	0.22
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	13	0.22
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	2	0.21
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	5	0.21
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	7	0.21
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	10	0.21
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	11	0.21
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	15	0.21
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	20	0.21
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	9	0.21
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	12	0.21
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	1	0.2
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	3	0.2
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	12	0.2
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	16	0.2
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	17	0.2
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	18	0.2
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	19	0.2
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	14	0.2
(2,45)	1:460:X:GLU:O	1:464:X:LYS:N	15	0.2
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	9	0.19
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	17	0.19
(2,48)	1:469:X:LYS:O	1:473:X:ALA:N	16	0.19
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	8	0.19
(2,53)	1:480:X:LEU:O	1:484:X:ASP:N	15	0.18
(2,50)	1:471:X:ARG:O	1:475:X:LEU:N	8	0.18
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	17	0.18
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	19	0.18
(1,764)	1:485:X:TYR:H	1:485:X:TYR:HD1	13	0.18
(1,764)	1:485:X:TYR:H	1:485:X:TYR:HD2	13	0.18
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	4	0.17

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	6	0.17
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	10	0.17
(2,50)	1:471:X:ARG:O	1:475:X:LEU:N	2	0.17
(2,50)	1:471:X:ARG:O	1:475:X:LEU:N	3	0.17
(2,50)	1:471:X:ARG:O	1:475:X:LEU:N	14	0.17
(2,50)	1:471:X:ARG:O	1:475:X:LEU:N	7	0.16
(2,50)	1:471:X:ARG:O	1:475:X:LEU:N	16	0.16
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	5	0.16
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	16	0.16
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	19	0.16
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	14	0.16
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	16	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	1	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	3	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	4	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	8	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	9	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	11	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	12	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	17	0.15
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	18	0.15
(1,764)	1:485:X:TYR:H	1:485:X:TYR:HD1	14	0.15
(1,764)	1:485:X:TYR:H	1:485:X:TYR:HD2	14	0.15
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	3	0.15
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	8	0.15
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	15	0.15
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	18	0.15
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	4	0.14
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	8	0.14
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	17	0.14
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	2	0.14
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	5	0.14
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	11	0.14
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	2	0.14
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	7	0.14
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	15	0.14
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	20	0.14
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	19	0.14
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	10	0.13
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	19	0.13
(2,51)	1:472:X:ALA:O	1:476:X:LEU:N	8	0.13
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	10	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1071)	1:489:X:LYS:H	1:489:X:LYS:HD3	5	0.13
(1,1059)	1:486:X:ARG:H	1:486:X:ARG:HG3	4	0.13
(1,1059)	1:486:X:ARG:H	1:486:X:ARG:HG3	6	0.13
(1,659)	1:472:X:ALA:H	1:474:X:LYS:HD3	6	0.13
(1,659)	1:472:X:ALA:H	1:474:X:LYS:HD3	19	0.13
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	7	0.13
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	9	0.13
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	12	0.13
(1,591)	1:450:X:ILE:HB	1:453:X:THR:H	16	0.13
(1,513)	1:451:X:GLU:H	1:453:X:THR:HG21	17	0.13
(1,513)	1:451:X:GLU:H	1:453:X:THR:HG22	17	0.13
(1,513)	1:451:X:GLU:H	1:453:X:THR:HG23	17	0.13
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	1	0.12
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	3	0.12
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	7	0.12
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	9	0.12
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	18	0.12
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	6	0.12
(1,1071)	1:489:X:LYS:H	1:489:X:LYS:HD3	19	0.12
(1,875)	1:496:X:LYS:HG3	1:497:X:PHE:H	12	0.12
(1,372)	1:458:X:ILE:HA	1:476:X:LEU:HG	13	0.12
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	13	0.11
(2,56)	1:483:X:LEU:O	1:487:X:ILE:N	14	0.11
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	2	0.11
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	5	0.11
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	11	0.11
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	12	0.11
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	20	0.11
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	10	0.11
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	20	0.11
(1,872)	1:496:X:LYS:HA	1:497:X:PHE:H	12	0.11
(1,863)	1:495:X:LYS:H	1:495:X:LYS:HG3	6	0.11
(1,372)	1:458:X:ILE:HA	1:476:X:LEU:HG	6	0.11
(1,342)	1:474:X:LYS:HA	1:474:X:LYS:HD3	6	0.11
(1,342)	1:474:X:LYS:HA	1:474:X:LYS:HD3	19	0.11
(1,246)	1:485:X:TYR:HD1	1:486:X:ARG:HD3	13	0.11
(1,246)	1:485:X:TYR:HD2	1:486:X:ARG:HD3	13	0.11
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	6	0.1
(2,52)	1:479:X:THR:O	1:483:X:LEU:N	14	0.1
(2,47)	1:468:X:VAL:O	1:472:X:ALA:N	1	0.1
(2,37)	1:482:X:GLN:O	1:486:X:ARG:H	13	0.1
(1,1059)	1:486:X:ARG:H	1:486:X:ARG:HG3	10	0.1

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

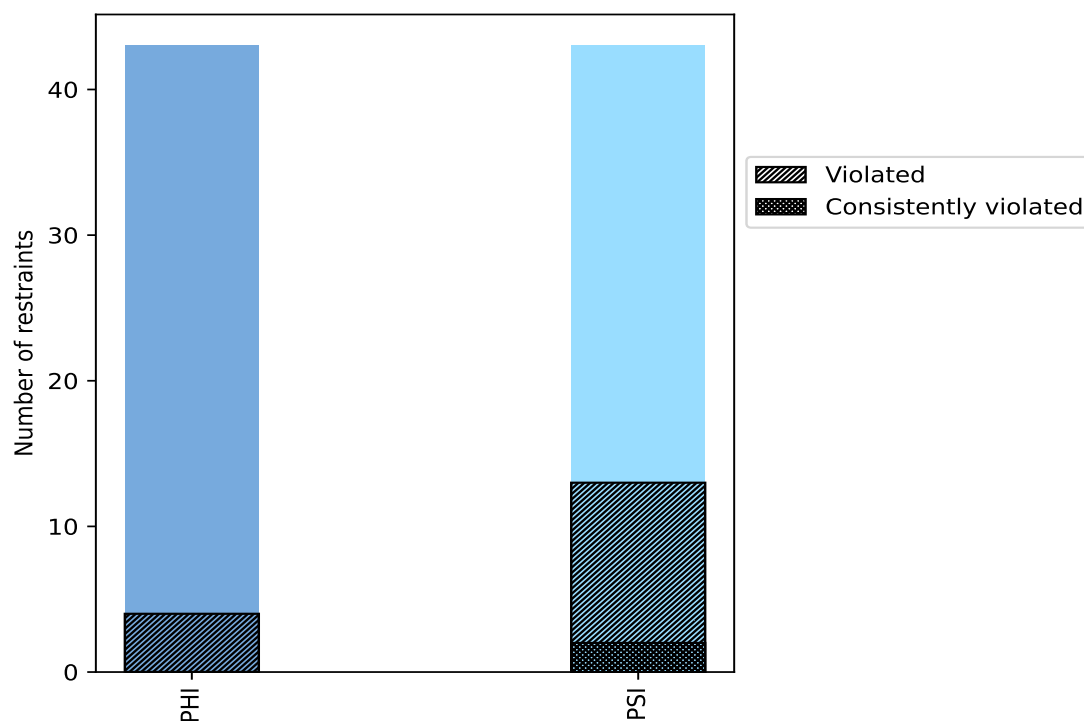
10.1 Summary of dihedral-angle violations [i](#)

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	43	50.0	4	9.3	4.7	0	0.0	0.0
PSI	43	50.0	13	30.2	15.1	2	4.7	2.3
Total	86	100.0	17	19.8	19.8	2	2.3	2.3

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](#)



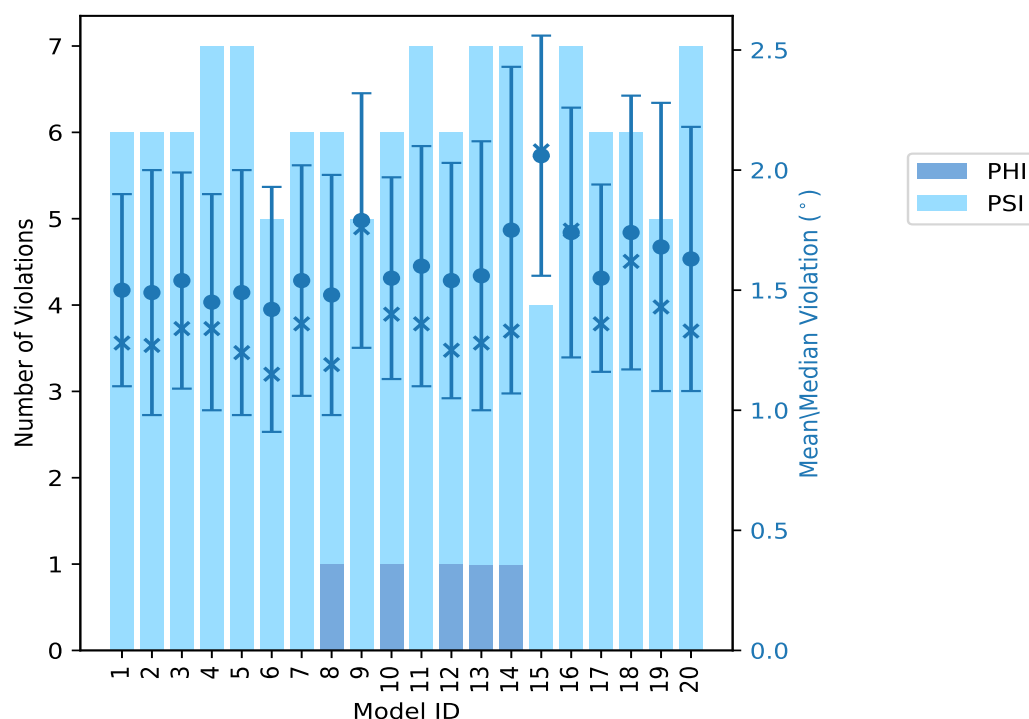
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

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	0	6	6	1.5	2.23	0.4	1.28
2	0	6	6	1.49	2.48	0.51	1.27
3	0	6	6	1.54	2.38	0.45	1.34
4	0	7	7	1.45	2.5	0.45	1.34
5	0	7	7	1.49	2.28	0.51	1.24
6	0	5	5	1.42	2.43	0.51	1.15
7	0	6	6	1.54	2.42	0.48	1.36
8	1	5	6	1.48	2.35	0.5	1.19
9	0	5	5	1.79	2.43	0.53	1.76
10	1	5	6	1.55	2.26	0.42	1.4
11	0	7	7	1.6	2.37	0.5	1.36
12	1	5	6	1.54	2.4	0.49	1.25
13	1	6	7	1.56	2.69	0.56	1.28
14	1	6	7	1.75	2.93	0.68	1.33
15	0	4	4	2.06	2.67	0.5	2.08
16	0	7	7	1.74	2.43	0.52	1.75
17	0	6	6	1.55	2.33	0.39	1.36
18	0	6	6	1.74	2.45	0.57	1.62
19	0	5	5	1.68	2.41	0.6	1.43
20	0	7	7	1.63	2.46	0.55	1.33

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

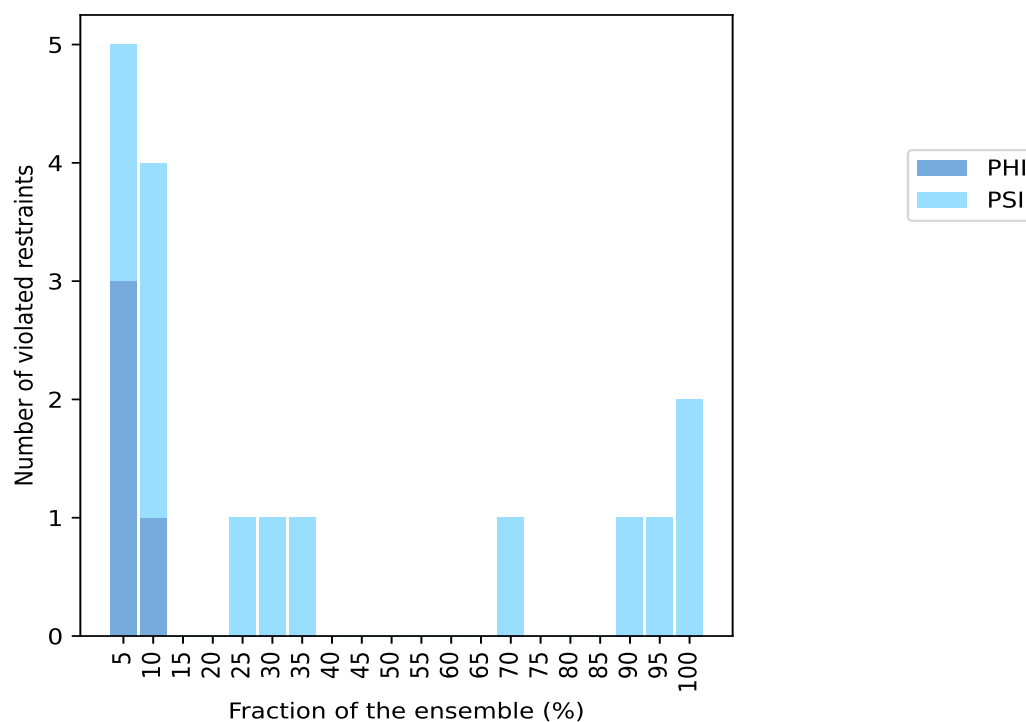
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Number of violated restraints			Fraction of the ensemble	
PHI	PSI	Total	Count ¹	%
0	0	0	12	60.0
0	0	0	13	65.0
0	1	1	14	70.0
0	0	0	15	75.0
0	0	0	16	80.0
0	0	0	17	85.0
0	1	1	18	90.0
0	1	1	19	95.0
0	2	2	20	100.0

¹ Number of models with violations

10.3.1 Bar graph : Dihedral-angle Violation statistics for the ensemble ⓘ

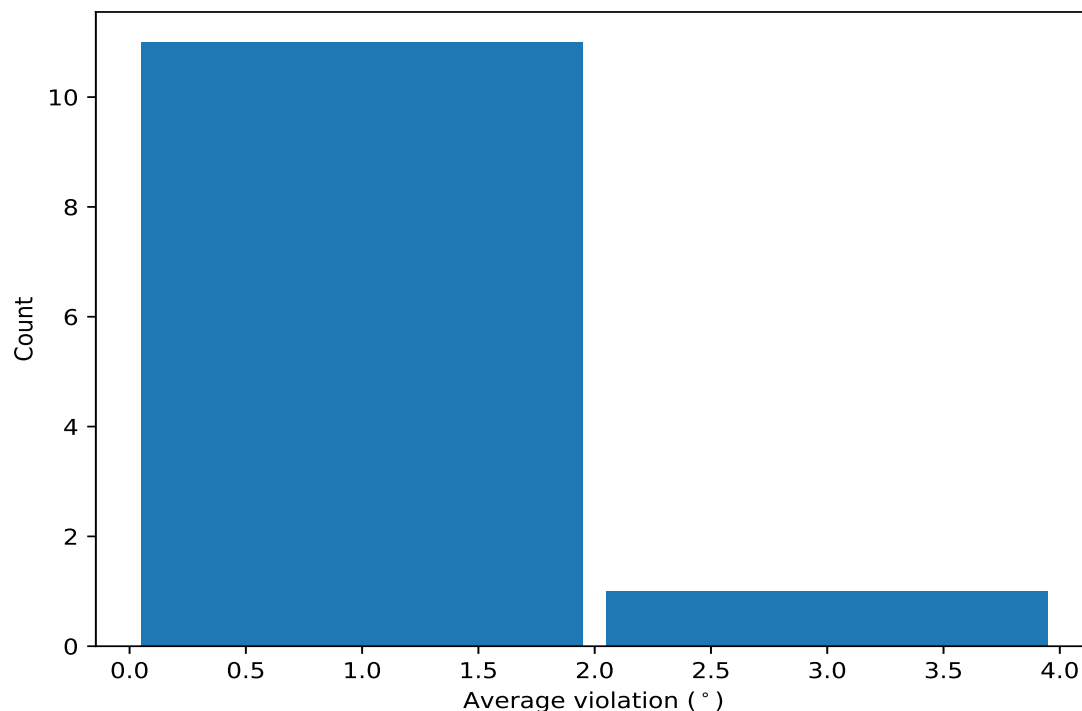


10.4 Most violated dihedral-angle restraints in the ensemble ⓘ

10.4.1 Histogram : Distribution of mean dihedral-angle violations ⓘ

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.

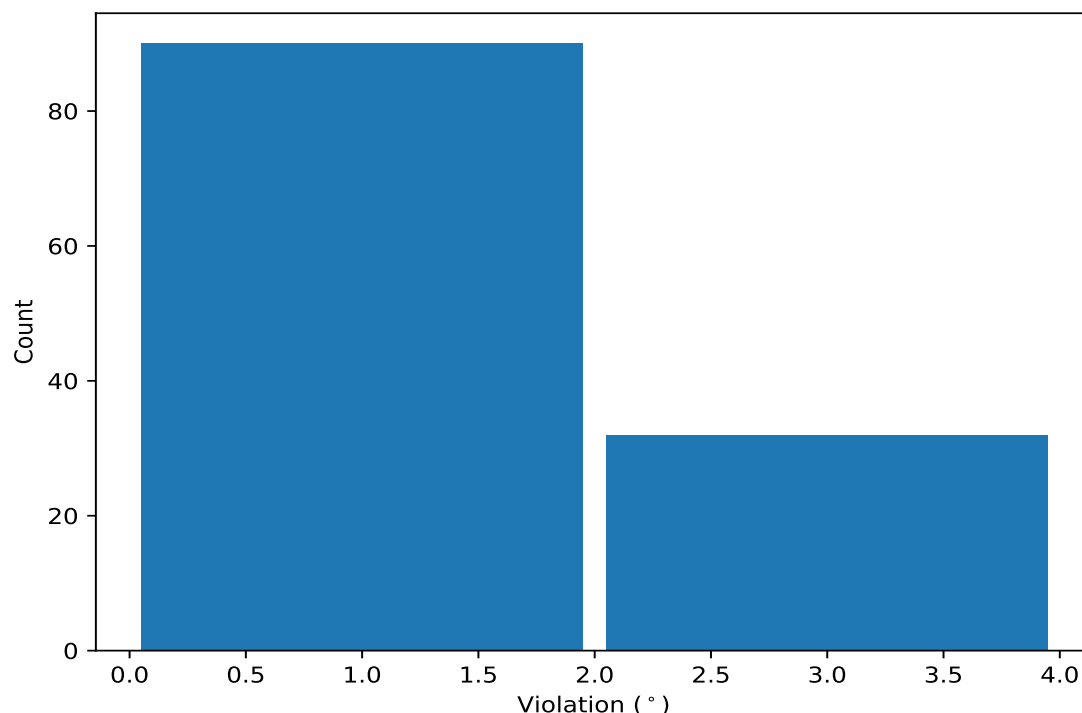
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models ¹	Mean	SD ²	Median
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	20	2.32	0.16	2.36
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	20	1.93	0.47	1.85
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	19	1.54	0.59	1.17
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	18	1.34	0.25	1.27
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	14	1.19	0.12	1.16
(1,54)	1:474:X:LYS:N	1:474:X:LYS:CA	1:474:X:LYS:C	1:475:X:LEU:N	7	1.29	0.04	1.29
(1,56)	1:475:X:LEU:N	1:475:X:LEU:CA	1:475:X:LEU:C	1:476:X:LEU:N	6	1.35	0.17	1.4
(1,20)	1:453:X:THR:N	1:453:X:THR:CA	1:453:X:THR:C	1:454:X:GLU:N	5	1.05	0.03	1.04
(1,69)	1:482:X:GLN:C	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	2	1.68	0.55	1.68
(1,4)	1:435:X:LEU:N	1:435:X:LEU:CA	1:435:X:LEU:C	1:436:X:LEU:N	2	1.28	0.04	1.28
(1,46)	1:470:X:SER:N	1:470:X:SER:CA	1:470:X:SER:C	1:471:X:ARG:N	2	1.25	0.08	1.25
(1,66)	1:481:X:ARG:N	1:481:X:ARG:CA	1:481:X:ARG:C	1:482:X:GLN:N	2	1.25	0.04	1.25

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 (°)
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	14	2.93
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	13	2.69
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	15	2.67
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	4	2.5
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	2	2.48
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	20	2.46
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	18	2.45
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	6	2.43
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	16	2.43
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	9	2.43
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	18	2.43
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	7	2.42
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	19	2.41
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	12	2.4

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	15	2.4
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	16	2.39
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	3	2.38
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	19	2.37
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	20	2.37
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	11	2.37
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	8	2.35
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	17	2.33
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	9	2.32
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	14	2.32
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	5	2.28
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	5	2.28
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	10	2.26
(1,69)	1:482:X:GLN:C	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	14	2.23
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	1	2.23
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	11	2.23
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	13	2.14
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	12	2.01
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	18	1.97
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	8	1.97
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	16	1.96
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	10	1.93
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	7	1.9
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	1	1.85
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	3	1.85
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	20	1.82
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	11	1.81
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	2	1.78
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	9	1.76
(1,38)	1:462:X:LEU:N	1:462:X:LEU:CA	1:462:X:LEU:C	1:463:X:GLU:N	15	1.76
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	16	1.75
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	17	1.75
(1,56)	1:475:X:LEU:N	1:475:X:LEU:CA	1:475:X:LEU:C	1:476:X:LEU:N	16	1.57
(1,3)	1:434:X:GLU:C	1:435:X:LEU:N	1:435:X:LEU:CA	1:435:X:LEU:C	10	1.56
(1,56)	1:475:X:LEU:N	1:475:X:LEU:CA	1:475:X:LEU:C	1:476:X:LEU:N	3	1.46
(1,56)	1:475:X:LEU:N	1:475:X:LEU:CA	1:475:X:LEU:C	1:476:X:LEU:N	7	1.46
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	4	1.45
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	19	1.43
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	4	1.42
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	15	1.41
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	17	1.39
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	11	1.36
(1,54)	1:474:X:LYS:N	1:474:X:LYS:CA	1:474:X:LYS:C	1:475:X:LEU:N	5	1.35
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	9	1.34
(1,56)	1:475:X:LEU:N	1:475:X:LEU:CA	1:475:X:LEU:C	1:476:X:LEU:N	2	1.34
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	4	1.34
(1,54)	1:474:X:LYS:N	1:474:X:LYS:CA	1:474:X:LYS:C	1:475:X:LEU:N	20	1.33
(1,46)	1:470:X:SER:N	1:470:X:SER:CA	1:470:X:SER:C	1:471:X:ARG:N	14	1.33
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	6	1.32
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	17	1.32
(1,4)	1:435:X:LEU:N	1:435:X:LEU:CA	1:435:X:LEU:C	1:436:X:LEU:N	17	1.32

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	20	1.31
(1,54)	1:474:X:LYS:N	1:474:X:LYS:CA	1:474:X:LYS:C	1:475:X:LEU:N	1	1.3
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	13	1.3
(1,54)	1:474:X:LYS:N	1:474:X:LYS:CA	1:474:X:LYS:C	1:475:X:LEU:N	11	1.29
(1,66)	1:481:X:ARG:N	1:481:X:ARG:CA	1:481:X:ARG:C	1:482:X:GLN:N	13	1.28
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	1	1.27
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	18	1.27
(1,54)	1:474:X:LYS:N	1:474:X:LYS:CA	1:474:X:LYS:C	1:475:X:LEU:N	18	1.27
(1,12)	1:449:X:PHE:N	1:449:X:PHE:CA	1:449:X:PHE:C	1:450:X:ILE:N	12	1.26
(1,54)	1:474:X:LYS:N	1:474:X:LYS:CA	1:474:X:LYS:C	1:475:X:LEU:N	10	1.25
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	7	1.25
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	5	1.24
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	12	1.24
(1,4)	1:435:X:LEU:N	1:435:X:LEU:CA	1:435:X:LEU:C	1:436:X:LEU:N	10	1.24
(1,71)	1:483:X:LEU:C	1:484:X:ASP:N	1:484:X:ASP:CA	1:484:X:ASP:C	8	1.22
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	3	1.22
(1,54)	1:474:X:LYS:N	1:474:X:LYS:CA	1:474:X:LYS:C	1:475:X:LEU:N	4	1.22
(1,66)	1:481:X:ARG:N	1:481:X:ARG:CA	1:481:X:ARG:C	1:482:X:GLN:N	14	1.21
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	7	1.2
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	3	1.2
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	2	1.19
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	13	1.18
(1,56)	1:475:X:LEU:N	1:475:X:LEU:CA	1:475:X:LEU:C	1:476:X:LEU:N	13	1.17
(1,46)	1:470:X:SER:N	1:470:X:SER:CA	1:470:X:SER:C	1:471:X:ARG:N	8	1.17
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	1	1.17
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	1	1.17
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	17	1.17
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	5	1.16
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	12	1.16
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	14	1.16
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	6	1.15
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	4	1.15
(1,69)	1:482:X:GLN:C	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	13	1.14
(1,17)	1:451:X:GLU:C	1:452:X:GLU:N	1:452:X:GLU:CA	1:452:X:GLU:C	12	1.14
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	19	1.14
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	6	1.13
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	8	1.13
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	11	1.12
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	3	1.12
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	5	1.1
(1,20)	1:453:X:THR:N	1:453:X:THR:CA	1:453:X:THR:C	1:454:X:GLU:N	20	1.1
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	2	1.09
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	10	1.08
(1,56)	1:475:X:LEU:N	1:475:X:LEU:CA	1:475:X:LEU:C	1:476:X:LEU:N	4	1.08
(1,52)	1:473:X:ALA:N	1:473:X:ALA:CA	1:473:X:ALA:C	1:474:X:LYS:N	14	1.08
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	9	1.08
(1,20)	1:453:X:THR:N	1:453:X:THR:CA	1:453:X:THR:C	1:454:X:GLU:N	16	1.07
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	6	1.06
(1,68)	1:482:X:GLN:N	1:482:X:GLN:CA	1:482:X:GLN:C	1:483:X:LEU:N	20	1.05
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	16	1.04
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	18	1.04

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,20)	1:453:X:THR:N	1:453:X:THR:CA	1:453:X:THR:C	1:454:X:GLU:N	2	1.04
(1,70)	1:483:X:LEU:N	1:483:X:LEU:CA	1:483:X:LEU:C	1:484:X:ASP:N	8	1.03
(1,32)	1:459:X:ILE:N	1:459:X:ILE:CA	1:459:X:ILE:C	1:460:X:GLU:N	19	1.03
(1,20)	1:453:X:THR:N	1:453:X:THR:CA	1:453:X:THR:C	1:454:X:GLU:N	11	1.03
(1,20)	1:453:X:THR:N	1:453:X:THR:CA	1:453:X:THR:C	1:454:X:GLU:N	5	1.01
(1,2)	1:432:X:LEU:N	1:432:X:LEU:CA	1:432:X:LEU:C	1:433:X:PRO:N	7	1.01