



Full wwPDB NMR Structure Validation Report ⓘ

Mar 4, 2026 – 11:10 PM UTC

PDB ID : 2N95 / pdb_00002n95
BMRB ID : 25881
Title : NMR structure of yeast Hit1 protein zinc finger
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Deposited on : 2015-11-06

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

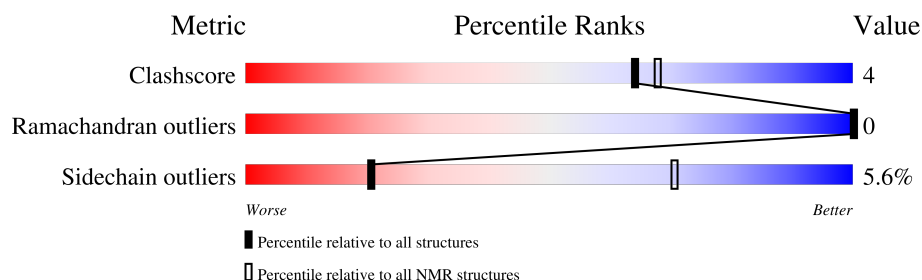
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 90%.

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	49	

2 Ensemble composition and analysis

This entry contains 20 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:6-A:37 (32)	0.19	7

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 2 single-model clusters were found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 747 atoms, of which 373 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Protein HIT1.

Mol	Chain	Residues	Atoms						Trace
1	A	49	Total	C	H	N	O	S	0
			745	229	373	70	66	7	

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P46973
A	-1	PRO	-	expression tag	UNP P46973
A	0	HIS	-	expression tag	UNP P46973

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

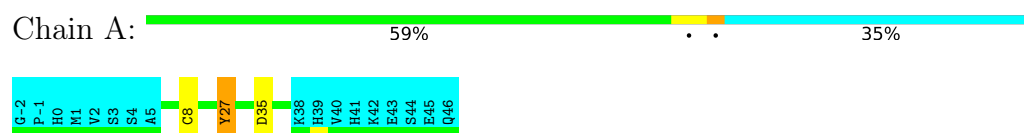
Mol	Chain	Residues	Atoms	
2	A	2	Total	Zn
			2	2

4 Residue-property plots [i](#)

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: Protein HIT1

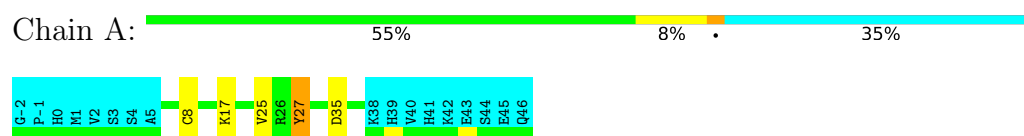


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

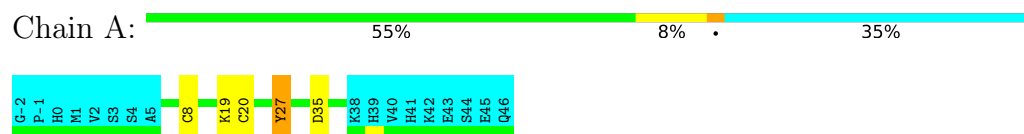
4.2.1 Score per residue for model 1

- Molecule 1: Protein HIT1



4.2.2 Score per residue for model 2

- Molecule 1: Protein HIT1



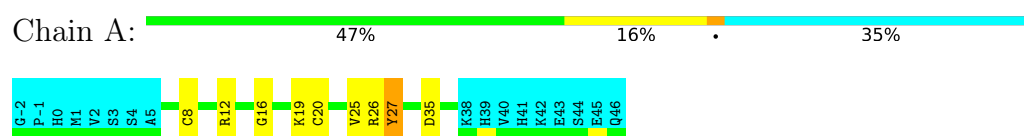
4.2.3 Score per residue for model 3

- Molecule 1: Protein HIT1



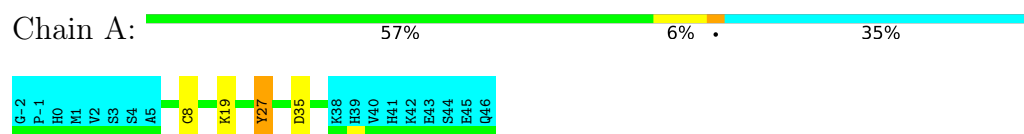
4.2.4 Score per residue for model 4

- Molecule 1: Protein HIT1



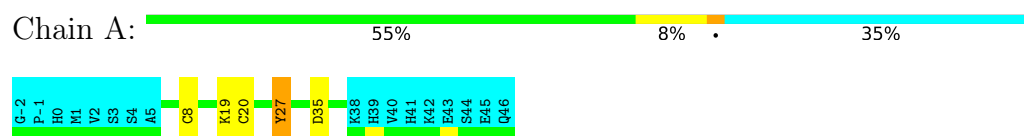
4.2.5 Score per residue for model 5

- Molecule 1: Protein HIT1



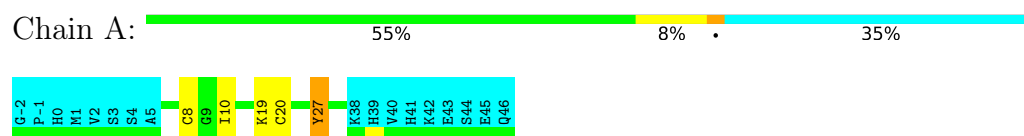
4.2.6 Score per residue for model 6

- Molecule 1: Protein HIT1



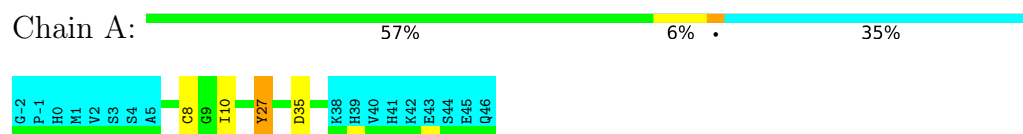
4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: Protein HIT1



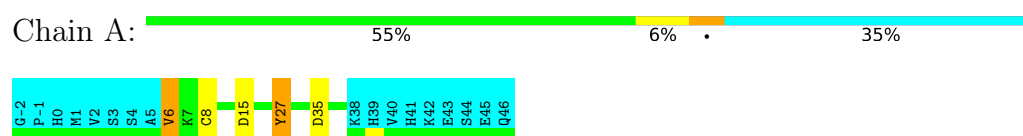
4.2.8 Score per residue for model 8

- Molecule 1: Protein HIT1



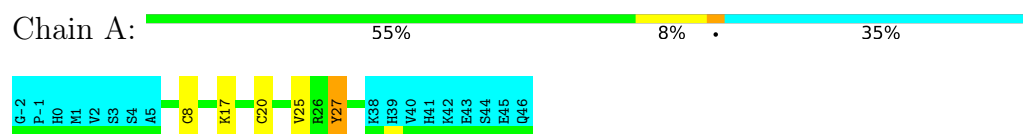
4.2.9 Score per residue for model 9

- Molecule 1: Protein HIT1



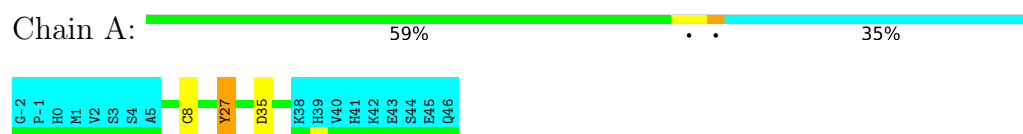
4.2.10 Score per residue for model 10

- Molecule 1: Protein HIT1



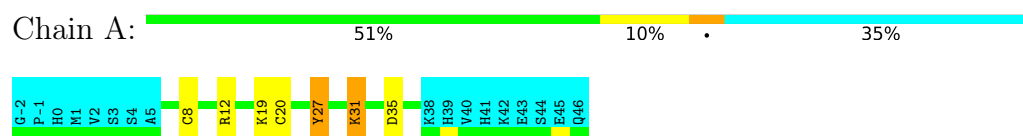
4.2.11 Score per residue for model 11

- Molecule 1: Protein HIT1



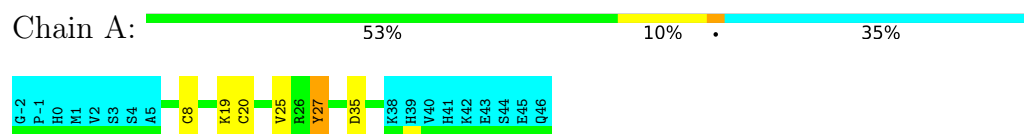
4.2.12 Score per residue for model 12

- Molecule 1: Protein HIT1



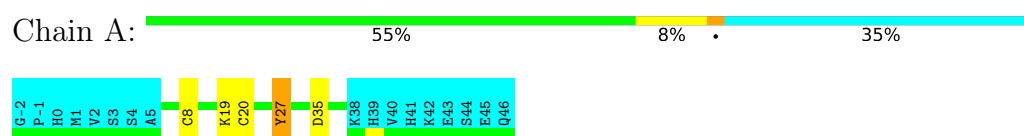
4.2.13 Score per residue for model 13

- Molecule 1: Protein HIT1



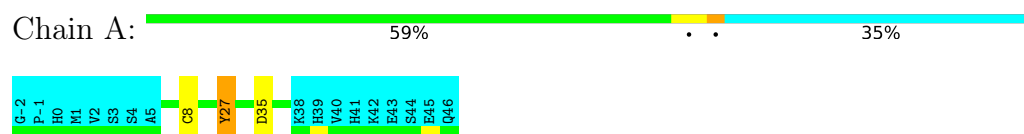
4.2.14 Score per residue for model 14

- Molecule 1: Protein HIT1



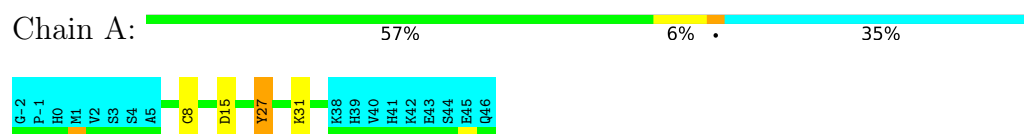
4.2.15 Score per residue for model 15

- Molecule 1: Protein HIT1



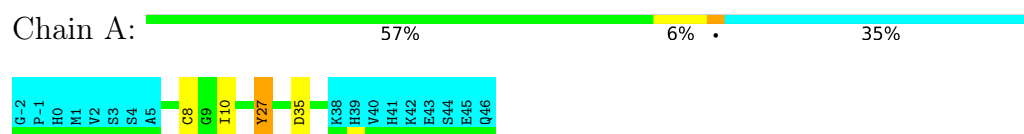
4.2.16 Score per residue for model 16

- Molecule 1: Protein HIT1



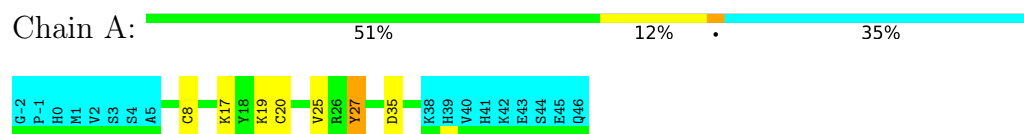
4.2.17 Score per residue for model 17

- Molecule 1: Protein HIT1



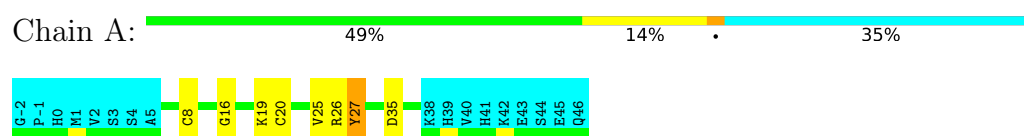
4.2.18 Score per residue for model 18

- Molecule 1: Protein HIT1



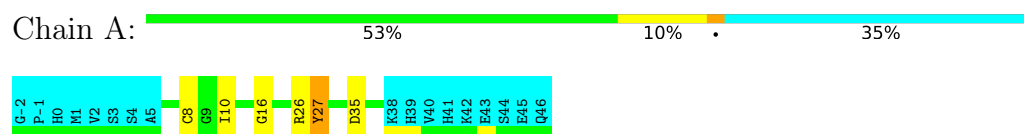
4.2.19 Score per residue for model 19

- Molecule 1: Protein HIT1



4.2.20 Score per residue for model 20

- Molecule 1: Protein HIT1



5 Refinement protocol and experimental data overview

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

Of the 250 calculated structures, 20 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	
CNS	structure solution	

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	962
Number of shifts mapped to atoms	962
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	90%

6 Model quality

6.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

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	A	240	246	246	2±1
All	All	4840	4920	4920	40

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:8:CYS:SG	1:A:27:TYR:HA	0.59	2.38	5	20
1:A:16:GLY:O	1:A:26:ARG:HD2	0.49	2.07	20	3
1:A:19:LYS:HD3	1:A:20:CYS:N	0.47	2.24	14	8
1:A:31:LYS:H	1:A:31:LYS:HD2	0.43	1.72	12	1
1:A:20:CYS:HB3	1:A:25:VAL:HG12	0.43	1.91	18	2
1:A:20:CYS:CB	1:A:25:VAL:HG12	0.42	2.45	10	3
1:A:6:VAL:O	1:A:15:ASP:HA	0.42	2.15	9	1
1:A:8:CYS:O	1:A:12:ARG:HD2	0.40	2.15	12	1
1:A:8:CYS:O	1:A:12:ARG:HA	0.40	2.17	4	1

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	32/49 (65%)	27±1 (85±4%)	5±1 (15±4%)	0±0 (0±0%)	100	100
All	All	640/980 (65%)	542 (85%)	98 (15%)	0 (0%)	100	100

There are no Ramachandran outliers.

6.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the sidechain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	26/41 (63%)	25±1 (94±2%)	1±1 (6±2%)	21	70
All	All	520/820 (63%)	491 (94%)	29 (6%)	21	70

All 6 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	27	TYR	20
1	A	17	LYS	3
1	A	19	LYS	2
1	A	31	LYS	2
1	A	6	VAL	1
1	A	15	ASP	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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 90% for the well-defined parts and 82% 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	520
Number of shifts mapped to atoms	520
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	47	-0.13 ± 0.37	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	41	0.27 ± 0.27	None needed (< 0.5 ppm)
$^{13}\text{C}'$	45	-0.12 ± 0.26	None needed (< 0.5 ppm)
^{15}N	42	-0.18 ± 0.75	None needed (< 0.5 ppm)

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 90%, i.e. 372 atoms were assigned a chemical shift out of a possible 414. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	161/162 (99%)	67/67 (100%)	64/64 (100%)	30/31 (97%)
Sidechain	199/225 (88%)	135/145 (93%)	64/68 (94%)	0/12 (0%)

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	Total	¹H	¹³C	¹⁵N
Aromatic	12/27 (44%)	6/12 (50%)	6/15 (40%)	0/0 (—%)
Overall	372/414 (90%)	208/224 (93%)	134/147 (91%)	30/43 (70%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 82%, i.e. 520 atoms were assigned a chemical shift out of a possible 633. 0 out of 6 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	225/246 (91%)	91/101 (90%)	92/98 (94%)	42/47 (89%)
Sidechain	271/336 (81%)	182/217 (84%)	88/104 (85%)	1/15 (7%)
Aromatic	24/51 (47%)	10/24 (42%)	10/21 (48%)	4/6 (67%)
Overall	520/633 (82%)	283/342 (83%)	190/223 (85%)	47/68 (69%)

7.1.4 Statistically unusual chemical shifts [i](#)

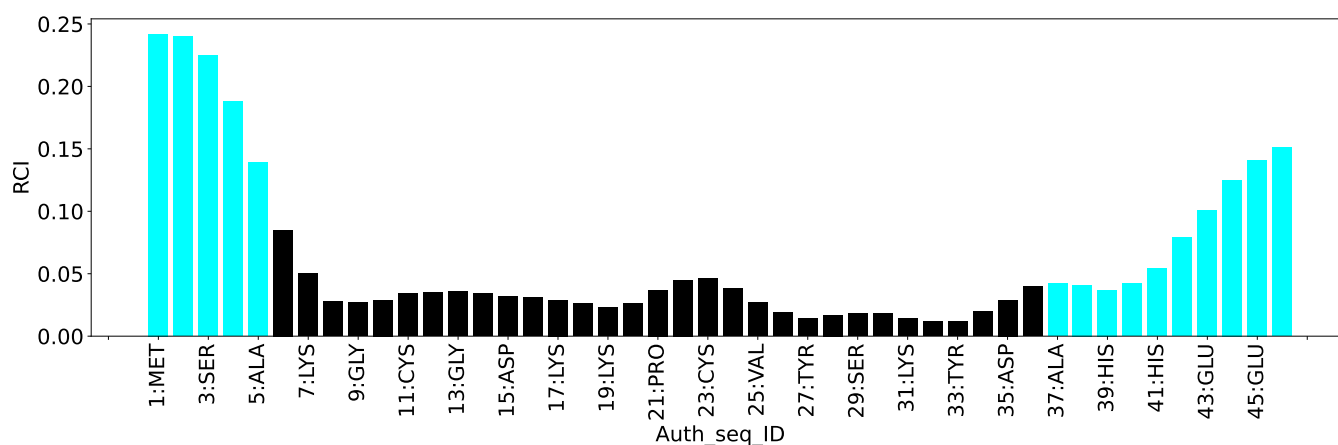
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	18	TYR	HB3	0.71	0.93 – 4.76	-5.6

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	442
Number of shifts mapped to atoms	442
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.2.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	43	-0.55 ± 0.37	None needed (imprecise)
$^{13}\text{C}_\beta$	38	0.22 ± 0.49	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	39	0.11 ± 0.41	None needed (< 0.5 ppm)

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 79%, i.e. 328 atoms were assigned a chemical shift out of a possible 414. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	127/162 (78%)	66/67 (99%)	31/64 (48%)	30/31 (97%)
Sidechain	189/225 (84%)	129/145 (89%)	60/68 (88%)	0/12 (0%)
Aromatic	12/27 (44%)	6/12 (50%)	6/15 (40%)	0/0 (—%)
Overall	328/414 (79%)	201/224 (90%)	97/147 (66%)	30/43 (70%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 70%, i.e. 442 atoms were assigned a chemical shift out of a possible 633. 0 out of 6 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	168/246 (68%)	86/101 (85%)	43/98 (44%)	39/47 (83%)
Sidechain	250/336 (74%)	170/217 (78%)	79/104 (76%)	1/15 (7%)
Aromatic	24/51 (47%)	10/24 (42%)	10/21 (48%)	4/6 (67%)
Overall	442/633 (70%)	266/342 (78%)	132/223 (59%)	44/68 (65%)

7.2.4 Statistically unusual chemical shifts [i](#)

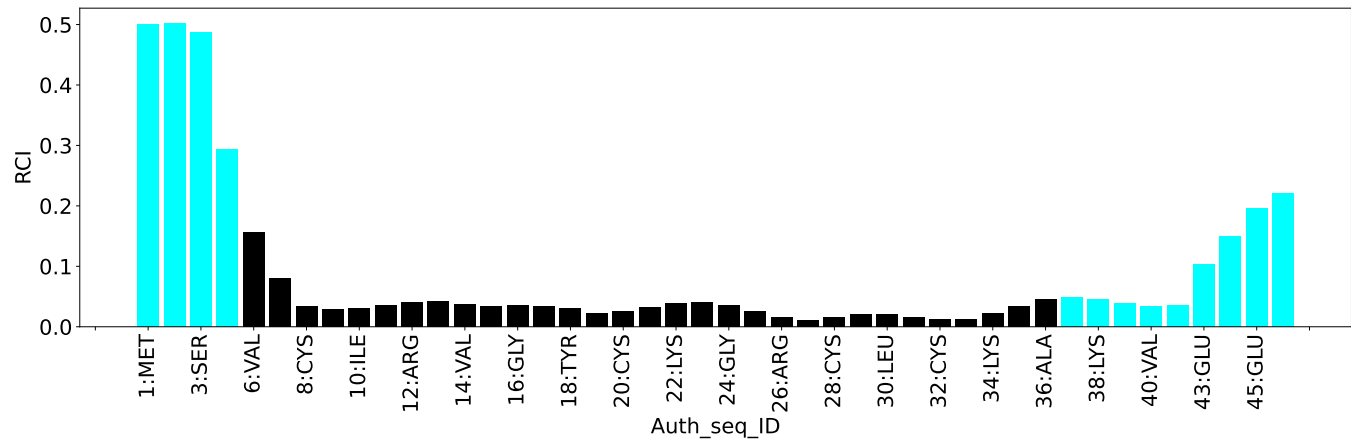
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
2	A	18	TYR	HB3	0.91	0.93 – 4.76	-5.0

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

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	958
Intra-residue ($ i-j =0$)	208
Sequential ($ i-j =1$)	266
Medium range ($ i-j >1$ and $ i-j <5$)	160
Long range ($ i-j \geq 5$)	310
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	10
Number of unmapped restraints	0
Number of restraints per residue	19.0
Number of long range restraints per residue ¹	6.1

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)	5.9	0.2
0.2-0.5 (Medium)	2.0	0.46
>0.5 (Large)	0.6	1.69

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)	1.1	5.4
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

9 Distance violation analysis

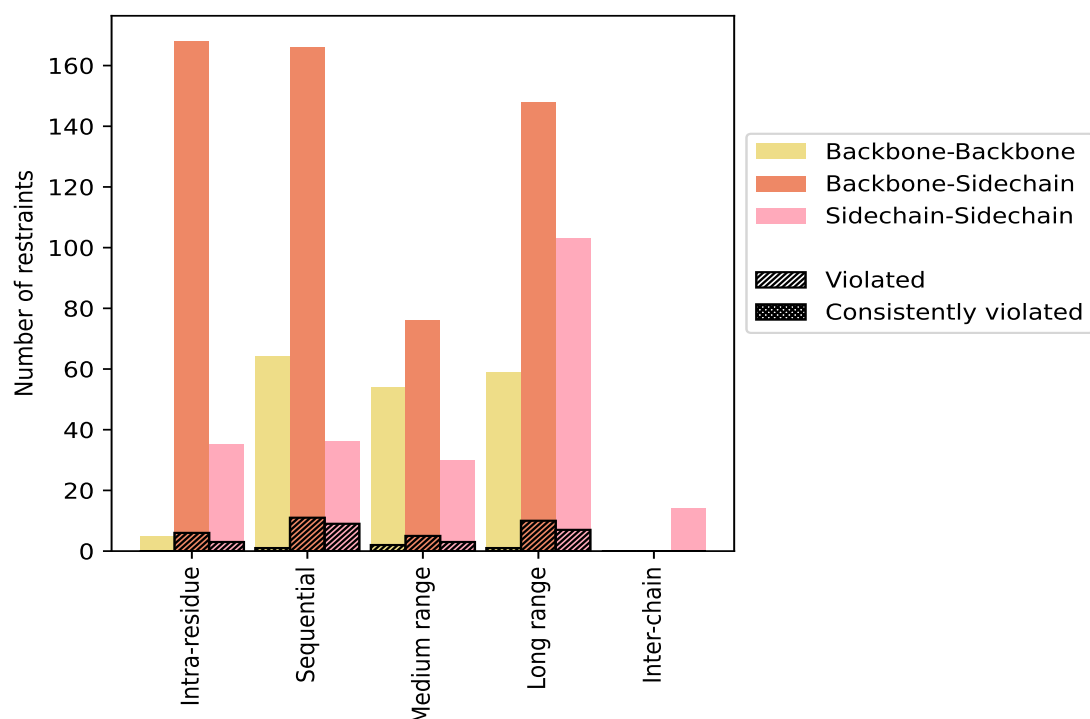
9.1 Summary of distance violations

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)	208	21.7	9	4.3	0.9	0	0.0	0.0
Backbone-Backbone	5	0.5	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	168	17.5	6	3.6	0.6	0	0.0	0.0
Sidechain-Sidechain	35	3.7	3	8.6	0.3	0	0.0	0.0
Sequential (i-j =1)	266	27.8	21	7.9	2.2	0	0.0	0.0
Backbone-Backbone	64	6.7	1	1.6	0.1	0	0.0	0.0
Backbone-Sidechain	166	17.3	11	6.6	1.1	0	0.0	0.0
Sidechain-Sidechain	36	3.8	9	25.0	0.9	0	0.0	0.0
Medium range (i-j >1 & i-j <5)	160	16.7	10	6.2	1.0	0	0.0	0.0
Backbone-Backbone	54	5.6	2	3.7	0.2	0	0.0	0.0
Backbone-Sidechain	76	7.9	5	6.6	0.5	0	0.0	0.0
Sidechain-Sidechain	30	3.1	3	10.0	0.3	0	0.0	0.0
Long range (i-j ≥5)	310	32.4	18	5.8	1.9	0	0.0	0.0
Backbone-Backbone	59	6.2	1	1.7	0.1	0	0.0	0.0
Backbone-Sidechain	148	15.4	10	6.8	1.0	0	0.0	0.0
Sidechain-Sidechain	103	10.8	7	6.8	0.7	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	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	958	100.0	58	6.1	6.1	0	0.0	0.0
Backbone-Backbone	182	19.0	4	2.2	0.4	0	0.0	0.0
Backbone-Sidechain	558	58.2	32	5.7	3.3	0	0.0	0.0
Sidechain-Sidechain	218	22.8	22	10.1	2.3	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	0	3	1	2	0	6	0.17	0.31	0.08	0.12
2	2	4	1	3	0	10	0.15	0.3	0.07	0.12
3	2	5	1	1	0	9	0.16	0.3	0.06	0.14
4	2	5	3	4	0	14	0.14	0.3	0.05	0.12
5	2	2	0	2	0	6	0.17	0.27	0.06	0.14
6	3	5	1	3	0	12	0.15	0.33	0.06	0.12
7	2	3	1	3	0	9	0.17	0.33	0.08	0.12
8	3	4	4	2	0	13	0.15	0.3	0.06	0.13
9	2	6	0	7	0	15	0.5	1.23	0.4	0.23
10	0	4	1	1	0	6	0.22	0.3	0.07	0.22

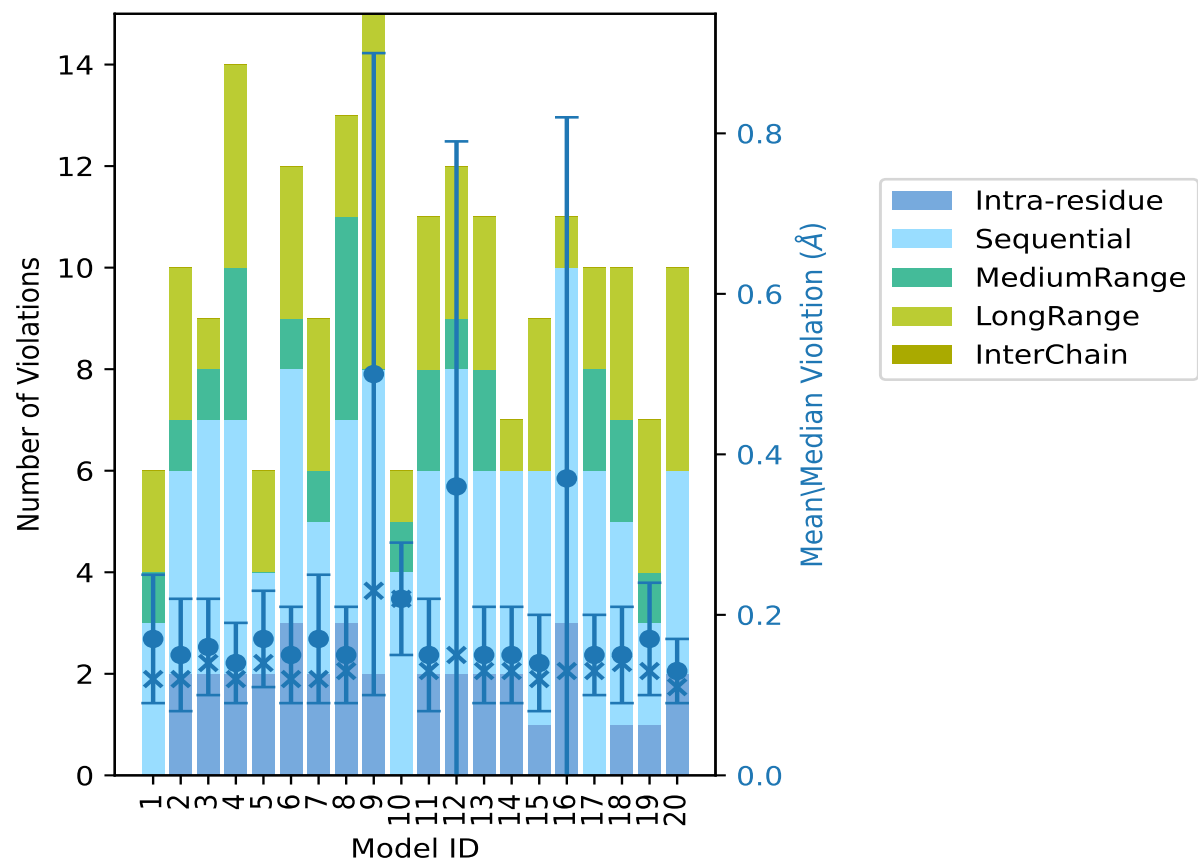
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
11	2	4	2	3	0	11	0.15	0.31	0.07	0.13
12	2	6	1	3	0	12	0.36	1.64	0.43	0.15
13	2	4	2	3	0	11	0.15	0.32	0.06	0.13
14	2	4	0	1	0	7	0.15	0.28	0.06	0.13
15	1	5	0	3	0	9	0.14	0.3	0.06	0.12
16	3	7	0	1	0	11	0.37	1.69	0.45	0.13
17	0	6	2	2	0	10	0.15	0.27	0.05	0.13
18	1	4	2	3	0	10	0.15	0.33	0.06	0.14
19	1	2	1	3	0	7	0.17	0.31	0.07	0.13
20	2	4	0	4	0	10	0.13	0.23	0.04	0.11

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

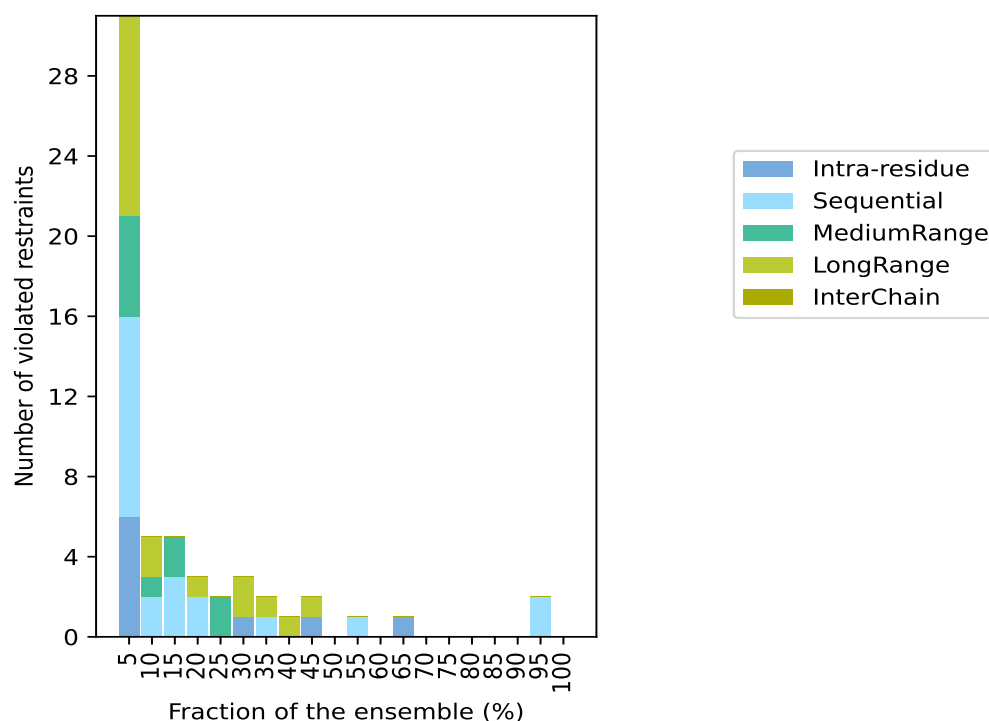
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 886(IR:199, SQ:245, MR:150, LR:292, IC:0) restraints are 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	%
6	10	5	10	0	31	1	5.0
0	2	1	2	0	5	2	10.0
0	3	2	0	0	5	3	15.0
0	2	0	1	0	3	4	20.0
0	0	2	0	0	2	5	25.0
1	0	0	2	0	3	6	30.0
0	1	0	1	0	2	7	35.0
0	0	0	1	0	1	8	40.0
1	0	0	1	0	2	9	45.0
0	0	0	0	0	0	10	50.0
0	1	0	0	0	1	11	55.0
0	0	0	0	0	0	12	60.0
1	0	0	0	0	1	13	65.0
0	0	0	0	0	0	14	70.0
0	0	0	0	0	0	15	75.0
0	0	0	0	0	0	16	80.0
0	0	0	0	0	0	17	85.0
0	0	0	0	0	0	18	90.0
0	2	0	0	0	2	19	95.0
0	0	0	0	0	0	20	100.0

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

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

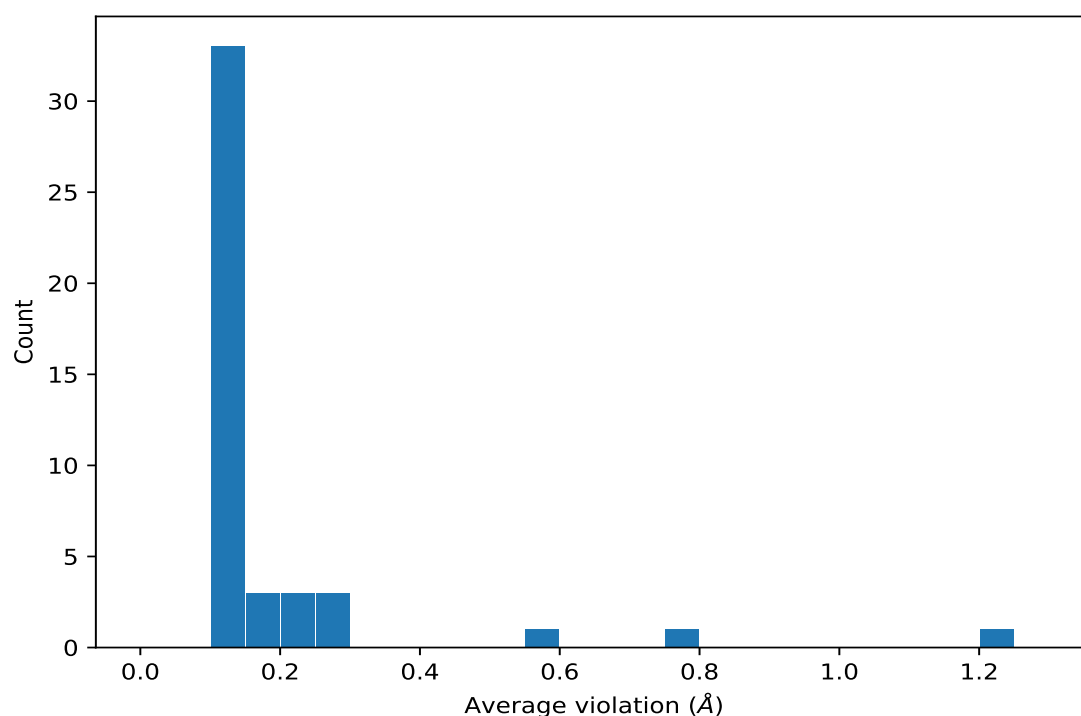
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



9.4 Most violated distance restraints in the ensemble [i](#)

9.4.1 Histogram : Distribution of mean distance violations [i](#)

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models in the ensemble



9.4.2 Table: Most violated distance restraints [i](#)

The following table provides the mean and the standard deviation of the violation for each restraint sorted by number of violated models and the mean value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	19	0.29	0.04	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	19	0.29	0.04	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	19	0.29	0.04	0.3
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	19	0.2	0.05	0.2
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	13	0.13	0.01	0.13
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	13	0.13	0.01	0.13
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	11	0.11	0.01	0.1
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	9	0.2	0.08	0.16
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	9	0.2	0.08	0.16
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	9	0.14	0.03	0.14
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	9	0.14	0.03	0.14
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	9	0.14	0.03	0.14
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	8	0.12	0.01	0.11
(1,816)	1:20:A:CYS:HB2	1:33:A:TYR:HE1	7	0.12	0.01	0.13
(1,590)	1:19:A:LYS:HG3	1:20:A:CYS:H	7	0.12	0.01	0.12
(1,15)	1:23:A:CYS:HB2	1:41:A:HIS:HD2	6	0.16	0.04	0.16

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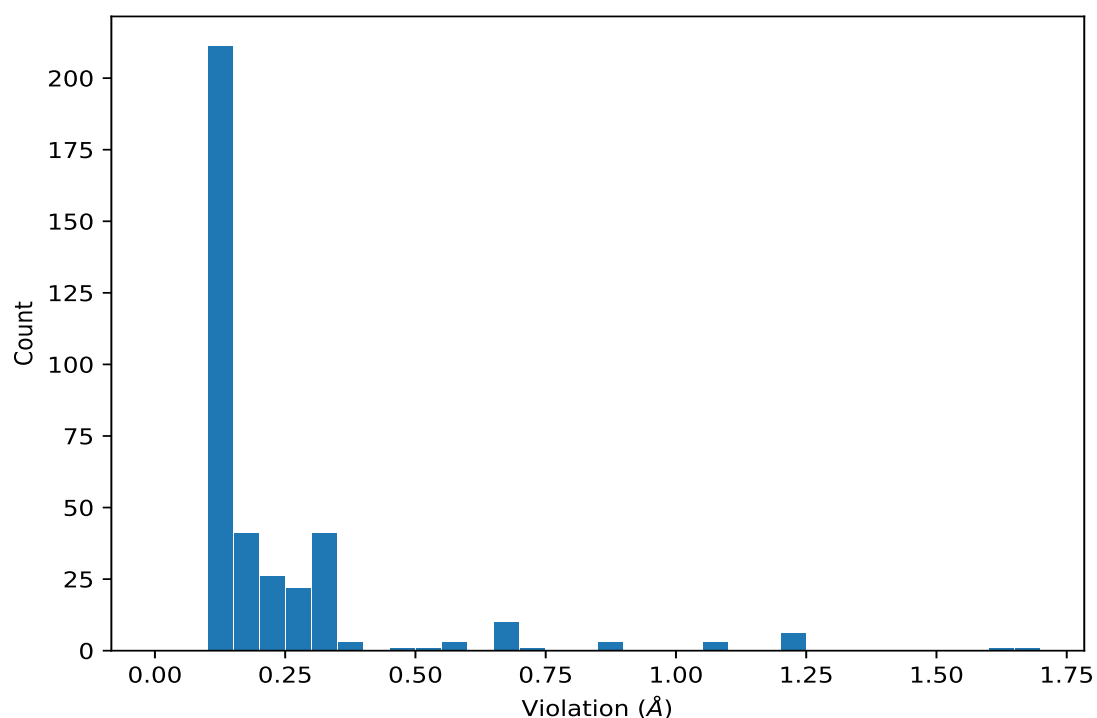
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE2	6	0.16	0.02	0.16
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE3	6	0.16	0.02	0.16
(1,182)	1:5:A:ALA:HB1	1:15:A:ASP:H	6	0.11	0.0	0.11
(1,182)	1:5:A:ALA:HB2	1:15:A:ASP:H	6	0.11	0.0	0.11
(1,182)	1:5:A:ALA:HB3	1:15:A:ASP:H	6	0.11	0.0	0.11
(1,508)	1:39:A:HIS:HE1	1:41:A:HIS:HD2	5	0.15	0.02	0.14
(1,45)	1:25:A:VAL:HG11	1:27:A:TYR:HE1	5	0.11	0.01	0.11
(1,45)	1:25:A:VAL:HG12	1:27:A:TYR:HE1	5	0.11	0.01	0.11
(1,45)	1:25:A:VAL:HG13	1:27:A:TYR:HE1	5	0.11	0.01	0.11
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG21	4	0.13	0.02	0.13
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG22	4	0.13	0.02	0.13
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG23	4	0.13	0.02	0.13
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG21	4	0.13	0.03	0.12
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG22	4	0.13	0.03	0.12
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG23	4	0.13	0.03	0.12
(1,798)	1:20:A:CYS:HB3	1:26:A:ARG:H	4	0.12	0.02	0.11
(1,459)	1:30:A:LEU:HB3	1:31:A:LYS:HG2	3	0.76	0.63	0.46
(1,248)	1:20:A:CYS:HB2	1:24:A:GLY:H	3	0.12	0.02	0.11
(1,401)	1:17:A:LYS:HD2	1:18:A:TYR:HB3	3	0.12	0.0	0.12
(1,401)	1:17:A:LYS:HD3	1:18:A:TYR:HB3	3	0.12	0.0	0.12
(1,764)	1:27:A:TYR:HD1	1:29:A:SER:HA	3	0.11	0.01	0.11
(1,302)	1:35:A:ASP:H	1:36:A:ALA:H	3	0.1	0.0	0.1
(1,460)	1:30:A:LEU:HB2	1:31:A:LYS:HG2	2	1.21	0.48	1.21
(1,287)	1:31:A:LYS:HG2	1:32:A:CYS:H	2	0.6	0.09	0.6
(1,197)	1:17:A:LYS:H	1:28:A:CYS:HB2	2	0.12	0.0	0.12
(1,641)	1:27:A:TYR:HB3	1:29:A:SER:H	2	0.12	0.0	0.12
(1,835)	1:14:A:VAL:HG11	1:28:A:CYS:HB2	2	0.1	0.0	0.1
(1,835)	1:14:A:VAL:HG12	1:28:A:CYS:HB2	2	0.1	0.0	0.1
(1,835)	1:14:A:VAL:HG13	1:28:A:CYS:HB2	2	0.1	0.0	0.1

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)
(1,460)	1:30:A:LEU:HB2	1:31:A:LYS:HG2	16	1.69
(1,459)	1:30:A:LEU:HB3	1:31:A:LYS:HG2	12	1.64
(1,399)	1:6:A:VAL:HG11	1:17:A:LYS:HA	9	1.23
(1,399)	1:6:A:VAL:HG12	1:17:A:LYS:HA	9	1.23
(1,399)	1:6:A:VAL:HG13	1:17:A:LYS:HA	9	1.23
(1,387)	1:6:A:VAL:HG11	1:15:A:ASP:HA	9	1.2
(1,387)	1:6:A:VAL:HG12	1:15:A:ASP:HA	9	1.2
(1,387)	1:6:A:VAL:HG13	1:15:A:ASP:HA	9	1.2
(1,718)	1:6:A:VAL:HG11	1:19:A:LYS:HA	9	1.05
(1,718)	1:6:A:VAL:HG12	1:19:A:LYS:HA	9	1.05
(1,718)	1:6:A:VAL:HG13	1:19:A:LYS:HA	9	1.05
(2,25)	1:6:A:VAL:HG11	1:27:A:TYR:HA	9	0.87
(2,25)	1:6:A:VAL:HG12	1:27:A:TYR:HA	9	0.87
(2,25)	1:6:A:VAL:HG13	1:27:A:TYR:HA	9	0.87
(1,460)	1:30:A:LEU:HB2	1:31:A:LYS:HG2	12	0.74
(1,346)	1:5:A:ALA:HB1	1:6:A:VAL:HG11	9	0.69

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,346)	1:5:A:ALA:HB1	1:6:A:VAL:HG12	9	0.69
(1,346)	1:5:A:ALA:HB1	1:6:A:VAL:HG13	9	0.69
(1,346)	1:5:A:ALA:HB2	1:6:A:VAL:HG11	9	0.69
(1,346)	1:5:A:ALA:HB2	1:6:A:VAL:HG12	9	0.69
(1,346)	1:5:A:ALA:HB2	1:6:A:VAL:HG13	9	0.69
(1,346)	1:5:A:ALA:HB3	1:6:A:VAL:HG11	9	0.69
(1,346)	1:5:A:ALA:HB3	1:6:A:VAL:HG12	9	0.69
(1,346)	1:5:A:ALA:HB3	1:6:A:VAL:HG13	9	0.69
(1,287)	1:31:A:LYS:HG2	1:32:A:CYS:H	16	0.68
(1,123)	1:6:A:VAL:H	1:6:A:VAL:HG11	9	0.58
(1,123)	1:6:A:VAL:H	1:6:A:VAL:HG12	9	0.58
(1,123)	1:6:A:VAL:H	1:6:A:VAL:HG13	9	0.58
(1,287)	1:31:A:LYS:HG2	1:32:A:CYS:H	12	0.51
(1,459)	1:30:A:LEU:HB3	1:31:A:LYS:HG2	16	0.46
(1,191)	1:6:A:VAL:HG11	1:16:A:GLY:H	9	0.38
(1,191)	1:6:A:VAL:HG12	1:16:A:GLY:H	9	0.38
(1,191)	1:6:A:VAL:HG13	1:16:A:GLY:H	9	0.38
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	6	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	6	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	6	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	7	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	7	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	7	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	18	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	18	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	18	0.33
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	13	0.32
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	13	0.32
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	13	0.32
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	16	0.32
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	16	0.32
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	16	0.32
(2,1)	1:33:A:TYR:HE1	1:34:A:LYS:HG3	11	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	1	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	1	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	1	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	12	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	12	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	12	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	19	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	19	0.31
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	19	0.31

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	2	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	2	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	2	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	4	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	4	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	4	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	10	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	10	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	10	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	15	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	15	0.3
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	15	0.3
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	3	0.3
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	3	0.3
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	8	0.3
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	8	0.3
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	16	0.29
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	14	0.28
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	14	0.28
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	14	0.28
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	10	0.27
(1,778)	1:30:A:LEU:HB3	1:31:A:LYS:HA	10	0.27
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	11	0.27
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	11	0.27
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	11	0.27
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	17	0.27
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	17	0.27
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	17	0.27
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	5	0.27
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	5	0.27
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	2	0.26
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	7	0.26
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	7	0.26
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	1	0.25
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	8	0.25
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	5	0.25
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	5	0.25
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	5	0.25
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	12	0.23
(1,440)	1:6:A:VAL:HG11	1:26:A:ARG:HD3	9	0.23
(1,440)	1:6:A:VAL:HG12	1:26:A:ARG:HD3	9	0.23
(1,440)	1:6:A:VAL:HG13	1:26:A:ARG:HD3	9	0.23

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	20	0.23
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	20	0.23
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	20	0.23
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	9	0.22
(1,15)	1:23:A:CYS:HB2	1:41:A:HIS:HD2	19	0.22
(2,9)	1:6:A:VAL:HG21	1:7:A:LYS:HB3	9	0.21
(2,9)	1:6:A:VAL:HG22	1:7:A:LYS:HB3	9	0.21
(2,9)	1:6:A:VAL:HG23	1:7:A:LYS:HB3	9	0.21
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	7	0.21
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	14	0.21
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	3	0.21
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	3	0.21
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	3	0.21
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG21	8	0.21
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG22	8	0.21
(1,344)	1:5:A:ALA:HA	1:6:A:VAL:HG23	8	0.21
(1,15)	1:23:A:CYS:HB2	1:41:A:HIS:HD2	6	0.21
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	3	0.2
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	4	0.2
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	19	0.2
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	19	0.2
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	19	0.2
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG21	3	0.19
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG22	3	0.19
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG23	3	0.19
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	17	0.19
(1,356)	1:6:A:VAL:HG21	1:7:A:LYS:HG2	9	0.19
(1,356)	1:6:A:VAL:HG21	1:7:A:LYS:HG3	9	0.19
(1,356)	1:6:A:VAL:HG22	1:7:A:LYS:HG2	9	0.19
(1,356)	1:6:A:VAL:HG22	1:7:A:LYS:HG3	9	0.19
(1,356)	1:6:A:VAL:HG23	1:7:A:LYS:HG2	9	0.19
(1,356)	1:6:A:VAL:HG23	1:7:A:LYS:HG3	9	0.19
(1,347)	1:6:A:VAL:HA	1:6:A:VAL:HG21	9	0.19
(1,347)	1:6:A:VAL:HA	1:6:A:VAL:HG22	9	0.19
(1,347)	1:6:A:VAL:HA	1:6:A:VAL:HG23	9	0.19
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE2	7	0.18
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE3	7	0.18
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	15	0.18
(1,562)	1:12:A:ARG:HD2	1:13:A:GLY:H	4	0.18
(1,562)	1:12:A:ARG:HD3	1:13:A:GLY:H	4	0.18
(1,508)	1:39:A:HIS:HE1	1:41:A:HIS:HD2	2	0.18
(1,459)	1:30:A:LEU:HB3	1:31:A:LYS:HG2	10	0.18

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE2	6	0.17
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE3	6	0.17
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	13	0.17
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	20	0.17
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG21	9	0.17
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG22	9	0.17
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG23	9	0.17
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	6	0.17
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	6	0.17
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	6	0.17
(1,15)	1:23:A:CYS:HB2	1:41:A:HIS:HD2	13	0.17
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE2	12	0.16
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE3	12	0.16
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	18	0.16
(1,508)	1:39:A:HIS:HE1	1:41:A:HIS:HD2	4	0.16
(1,295)	1:35:A:ASP:H	1:39:A:HIS:H	8	0.16
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	6	0.16
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	6	0.16
(1,15)	1:23:A:CYS:HB2	1:41:A:HIS:HD2	18	0.16
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE2	13	0.15
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE3	13	0.15
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE2	18	0.14
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE3	18	0.14
(1,816)	1:20:A:CYS:HB2	1:33:A:TYR:HE1	2	0.14
(1,798)	1:20:A:CYS:HB3	1:26:A:ARG:H	17	0.14
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	5	0.14
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	11	0.14
(1,590)	1:19:A:LYS:HG3	1:20:A:CYS:H	15	0.14
(1,508)	1:39:A:HIS:HE1	1:41:A:HIS:HD2	10	0.14
(1,508)	1:39:A:HIS:HE1	1:41:A:HIS:HD2	17	0.14
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	11	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	3	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	3	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	4	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	4	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	8	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	8	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	12	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	12	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	20	0.14
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	20	0.14
(1,248)	1:20:A:CYS:HB2	1:24:A:GLY:H	13	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,241)	1:22:A:LYS:HG2	1:23:A:CYS:H	2	0.14
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	4	0.14
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	4	0.14
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	4	0.14
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	11	0.14
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	11	0.14
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	11	0.14
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	18	0.14
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	18	0.14
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	18	0.14
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG21	8	0.13
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG22	8	0.13
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG23	8	0.13
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE2	14	0.13
(1,834)	1:19:A:LYS:HB2	1:19:A:LYS:HE3	14	0.13
(1,816)	1:20:A:CYS:HB2	1:33:A:TYR:HE1	1	0.13
(1,816)	1:20:A:CYS:HB2	1:33:A:TYR:HE1	10	0.13
(1,816)	1:20:A:CYS:HB2	1:33:A:TYR:HE1	13	0.13
(1,816)	1:20:A:CYS:HB2	1:33:A:TYR:HE1	18	0.13
(1,779)	1:40:A:VAL:HA	1:41:A:HIS:HD2	6	0.13
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG21	8	0.13
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG22	8	0.13
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG23	8	0.13
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG21	17	0.13
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG22	17	0.13
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG23	17	0.13
(1,590)	1:19:A:LYS:HG3	1:20:A:CYS:H	9	0.13
(1,590)	1:19:A:LYS:HG3	1:20:A:CYS:H	16	0.13
(1,508)	1:39:A:HIS:HE1	1:41:A:HIS:HD2	8	0.13
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	5	0.13
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	17	0.13
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	5	0.13
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	5	0.13
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	19	0.13
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	19	0.13
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	4	0.13
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	11	0.13
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	11	0.13
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	15	0.13
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	15	0.13
(1,46)	1:10:A:ILE:HD11	1:33:A:TYR:HD1	14	0.13
(1,46)	1:10:A:ILE:HD12	1:33:A:TYR:HD1	14	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,46)	1:10:A:ILE:HD13	1:33:A:TYR:HD1	14	0.13
(1,45)	1:25:A:VAL:HG11	1:27:A:TYR:HE1	13	0.13
(1,45)	1:25:A:VAL:HG12	1:27:A:TYR:HE1	13	0.13
(1,45)	1:25:A:VAL:HG13	1:27:A:TYR:HE1	13	0.13
(1,894)	1:39:A:HIS:H	1:40:A:VAL:HG11	17	0.12
(1,894)	1:39:A:HIS:H	1:40:A:VAL:HG12	17	0.12
(1,894)	1:39:A:HIS:H	1:40:A:VAL:HG13	17	0.12
(1,894)	1:39:A:HIS:H	1:40:A:VAL:HG21	17	0.12
(1,894)	1:39:A:HIS:H	1:40:A:VAL:HG22	17	0.12
(1,894)	1:39:A:HIS:H	1:40:A:VAL:HG23	17	0.12
(1,764)	1:27:A:TYR:HD1	1:29:A:SER:HA	12	0.12
(1,671)	1:44:A:SER:HA	1:46:A:GLN:H	19	0.12
(1,641)	1:27:A:TYR:HB3	1:29:A:SER:H	18	0.12
(1,590)	1:19:A:LYS:HG3	1:20:A:CYS:H	3	0.12
(1,590)	1:19:A:LYS:HG3	1:20:A:CYS:H	20	0.12
(1,401)	1:17:A:LYS:HD2	1:18:A:TYR:HB3	15	0.12
(1,401)	1:17:A:LYS:HD3	1:18:A:TYR:HB3	15	0.12
(1,401)	1:17:A:LYS:HD2	1:18:A:TYR:HB3	17	0.12
(1,401)	1:17:A:LYS:HD3	1:18:A:TYR:HB3	17	0.12
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	6	0.12
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	6	0.12
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	13	0.12
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	13	0.12
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	14	0.12
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	14	0.12
(1,197)	1:17:A:LYS:H	1:28:A:CYS:HB2	12	0.12
(1,182)	1:5:A:ALA:HB1	1:15:A:ASP:H	3	0.12
(1,182)	1:5:A:ALA:HB2	1:15:A:ASP:H	3	0.12
(1,182)	1:5:A:ALA:HB3	1:15:A:ASP:H	3	0.12
(1,182)	1:5:A:ALA:HB1	1:15:A:ASP:H	7	0.12
(1,182)	1:5:A:ALA:HB2	1:15:A:ASP:H	7	0.12
(1,182)	1:5:A:ALA:HB3	1:15:A:ASP:H	7	0.12
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	16	0.12
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	16	0.12
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	16	0.12
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	13	0.12
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	13	0.12
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	13	0.12
(1,45)	1:25:A:VAL:HG11	1:27:A:TYR:HE1	4	0.12
(1,45)	1:25:A:VAL:HG12	1:27:A:TYR:HE1	4	0.12
(1,45)	1:25:A:VAL:HG13	1:27:A:TYR:HE1	4	0.12
(1,15)	1:23:A:CYS:HB2	1:41:A:HIS:HD2	11	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,3)	1:39:A:HIS:HE1	1:41:A:HIS:HE1	17	0.12
(2,30)	1:27:A:TYR:HA	1:33:A:TYR:HA	1	0.11
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG21	20	0.11
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG22	20	0.11
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG23	20	0.11
(1,798)	1:20:A:CYS:HB3	1:26:A:ARG:H	2	0.11
(1,798)	1:20:A:CYS:HB3	1:26:A:ARG:H	8	0.11
(1,775)	1:31:A:LYS:HA	1:31:A:LYS:HE2	16	0.11
(1,775)	1:31:A:LYS:HA	1:31:A:LYS:HE3	16	0.11
(1,764)	1:27:A:TYR:HD1	1:29:A:SER:HA	8	0.11
(1,641)	1:27:A:TYR:HB3	1:29:A:SER:H	4	0.11
(1,590)	1:19:A:LYS:HG3	1:20:A:CYS:H	18	0.11
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	9	0.11
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	19	0.11
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	20	0.11
(1,401)	1:17:A:LYS:HD2	1:18:A:TYR:HB3	7	0.11
(1,401)	1:17:A:LYS:HD3	1:18:A:TYR:HB3	7	0.11
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	2	0.11
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	2	0.11
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	11	0.11
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	11	0.11
(1,353)	1:7:A:LYS:HB2	1:7:A:LYS:HD2	4	0.11
(1,353)	1:7:A:LYS:HB2	1:7:A:LYS:HD3	4	0.11
(1,302)	1:35:A:ASP:H	1:36:A:ALA:H	6	0.11
(1,273)	1:29:A:SER:HB2	1:31:A:LYS:H	11	0.11
(1,248)	1:20:A:CYS:HB2	1:24:A:GLY:H	18	0.11
(1,197)	1:17:A:LYS:H	1:28:A:CYS:HB2	7	0.11
(1,182)	1:5:A:ALA:HB1	1:15:A:ASP:H	4	0.11
(1,182)	1:5:A:ALA:HB2	1:15:A:ASP:H	4	0.11
(1,182)	1:5:A:ALA:HB3	1:15:A:ASP:H	4	0.11
(1,182)	1:5:A:ALA:HB1	1:15:A:ASP:H	8	0.11
(1,182)	1:5:A:ALA:HB2	1:15:A:ASP:H	8	0.11
(1,182)	1:5:A:ALA:HB3	1:15:A:ASP:H	8	0.11
(1,182)	1:5:A:ALA:HB1	1:15:A:ASP:H	15	0.11
(1,182)	1:5:A:ALA:HB2	1:15:A:ASP:H	15	0.11
(1,182)	1:5:A:ALA:HB3	1:15:A:ASP:H	15	0.11
(1,182)	1:5:A:ALA:HB1	1:15:A:ASP:H	20	0.11
(1,182)	1:5:A:ALA:HB2	1:15:A:ASP:H	20	0.11
(1,182)	1:5:A:ALA:HB3	1:15:A:ASP:H	20	0.11
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	1	0.11
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	13	0.11
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	15	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,106)	1:33:A:TYR:HD1	1:34:A:LYS:HG3	11	0.11
(1,68)	1:33:A:TYR:HD1	1:34:A:LYS:HG2	4	0.11
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	2	0.11
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	2	0.11
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	2	0.11
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	5	0.11
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	5	0.11
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	5	0.11
(1,48)	1:25:A:VAL:HG11	1:41:A:HIS:HD2	20	0.11
(1,48)	1:25:A:VAL:HG12	1:41:A:HIS:HD2	20	0.11
(1,48)	1:25:A:VAL:HG13	1:41:A:HIS:HD2	20	0.11
(1,45)	1:25:A:VAL:HG11	1:27:A:TYR:HE1	11	0.11
(1,45)	1:25:A:VAL:HG12	1:27:A:TYR:HE1	11	0.11
(1,45)	1:25:A:VAL:HG13	1:27:A:TYR:HE1	11	0.11
(1,5)	1:19:A:LYS:HA	1:27:A:TYR:HE1	12	0.11
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG21	6	0.1
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG22	6	0.1
(2,4)	1:5:A:ALA:H	1:6:A:VAL:HG23	6	0.1
(1,835)	1:14:A:VAL:HG11	1:28:A:CYS:HB2	12	0.1
(1,835)	1:14:A:VAL:HG12	1:28:A:CYS:HB2	12	0.1
(1,835)	1:14:A:VAL:HG13	1:28:A:CYS:HB2	12	0.1
(1,835)	1:14:A:VAL:HG11	1:28:A:CYS:HB2	15	0.1
(1,835)	1:14:A:VAL:HG12	1:28:A:CYS:HB2	15	0.1
(1,835)	1:14:A:VAL:HG13	1:28:A:CYS:HB2	15	0.1
(1,816)	1:20:A:CYS:HB2	1:33:A:TYR:HE1	4	0.1
(1,816)	1:20:A:CYS:HB2	1:33:A:TYR:HE1	6	0.1
(1,798)	1:20:A:CYS:HB3	1:26:A:ARG:H	16	0.1
(1,764)	1:27:A:TYR:HD1	1:29:A:SER:HA	7	0.1
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG21	14	0.1
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG22	14	0.1
(1,603)	1:24:A:GLY:H	1:25:A:VAL:HG23	14	0.1
(1,590)	1:19:A:LYS:HG3	1:20:A:CYS:H	14	0.1
(1,482)	1:38:A:LYS:HA	1:38:A:LYS:HG3	20	0.1
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	7	0.1
(1,414)	1:20:A:CYS:HA	1:33:A:TYR:HE1	15	0.1
(1,412)	1:10:A:ILE:HG21	1:38:A:LYS:HE2	20	0.1
(1,412)	1:10:A:ILE:HG21	1:38:A:LYS:HE3	20	0.1
(1,412)	1:10:A:ILE:HG22	1:38:A:LYS:HE2	20	0.1
(1,412)	1:10:A:ILE:HG22	1:38:A:LYS:HE3	20	0.1
(1,412)	1:10:A:ILE:HG23	1:38:A:LYS:HE2	20	0.1
(1,412)	1:10:A:ILE:HG23	1:38:A:LYS:HE3	20	0.1
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD2	16	0.1

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,397)	1:17:A:LYS:HA	1:17:A:LYS:HD3	16	0.1
(1,302)	1:35:A:ASP:H	1:36:A:ALA:H	13	0.1
(1,302)	1:35:A:ASP:H	1:36:A:ALA:H	18	0.1
(1,274)	1:29:A:SER:HB3	1:31:A:LYS:H	8	0.1
(1,248)	1:20:A:CYS:HB2	1:24:A:GLY:H	6	0.1
(1,163)	1:12:A:ARG:H	1:12:A:ARG:HG2	8	0.1
(1,163)	1:12:A:ARG:H	1:12:A:ARG:HG3	8	0.1
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	2	0.1
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	3	0.1
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	6	0.1
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	12	0.1
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	17	0.1
(1,148)	1:10:A:ILE:H	1:11:A:CYS:HB3	19	0.1
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE2	2	0.1
(1,94)	1:7:A:LYS:HB3	1:7:A:LYS:HE3	2	0.1
(1,45)	1:25:A:VAL:HG11	1:27:A:TYR:HE1	1	0.1
(1,45)	1:25:A:VAL:HG12	1:27:A:TYR:HE1	1	0.1
(1,45)	1:25:A:VAL:HG13	1:27:A:TYR:HE1	1	0.1
(1,45)	1:25:A:VAL:HG11	1:27:A:TYR:HE1	3	0.1
(1,45)	1:25:A:VAL:HG12	1:27:A:TYR:HE1	3	0.1
(1,45)	1:25:A:VAL:HG13	1:27:A:TYR:HE1	3	0.1
(1,15)	1:23:A:CYS:HB2	1:41:A:HIS:HD2	4	0.1

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

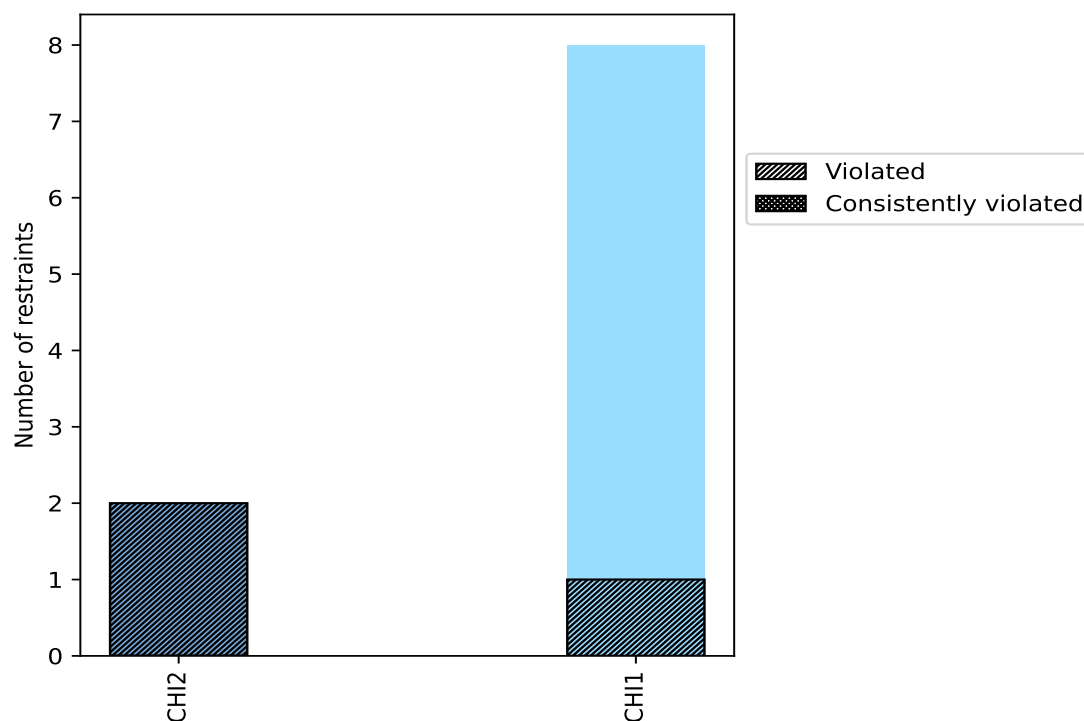
10.1 Summary of dihedral-angle violations [i](#)

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
CHI2	2	20.0	2	100.0	20.0	0	0.0	0.0
CHI1	8	80.0	1	12.5	10.0	0	0.0	0.0
Total	10	100.0	3	30.0	30.0	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](#)



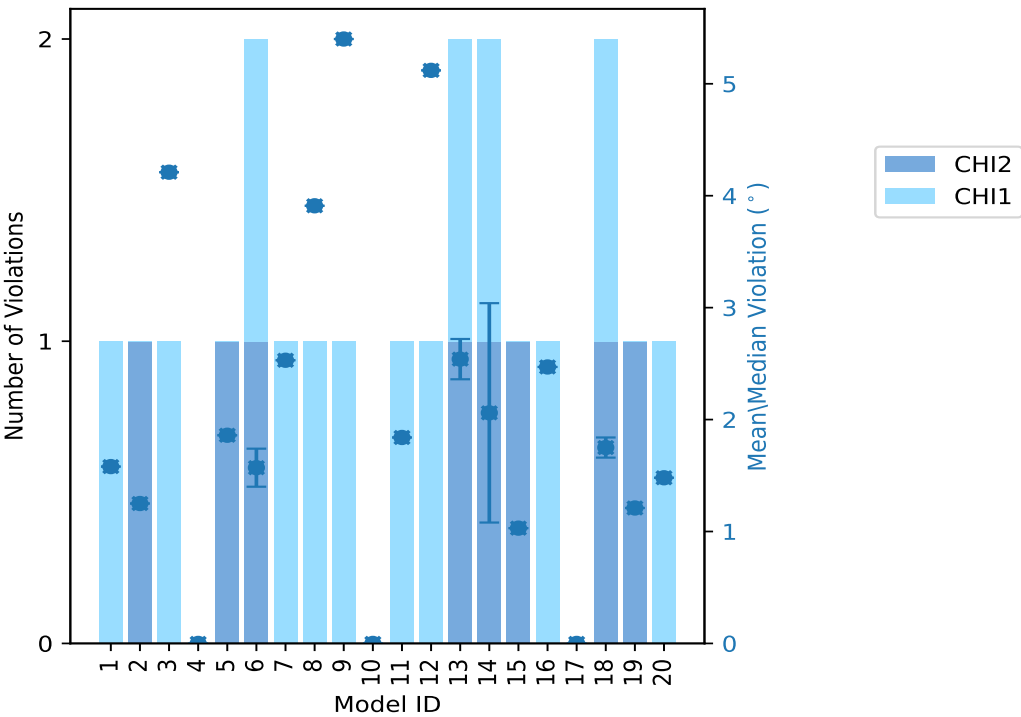
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

10.2 Dihedral-angle violation statistics for each model

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 (°)
	CHI2	CHI1	Total				
1	0	1	1	1.58	1.58	0.0	1.58
2	1	0	1	1.25	1.25	0.0	1.25
3	0	1	1	4.21	4.21	0.0	4.21
4	0	0	0	0.0	0.0	0.0	0.0
5	1	0	1	1.86	1.86	0.0	1.86
6	1	1	2	1.57	1.74	0.17	1.57
7	0	1	1	2.53	2.53	0.0	2.53
8	0	1	1	3.91	3.91	0.0	3.91
9	0	1	1	5.4	5.4	0.0	5.4
10	0	0	0	0.0	0.0	0.0	0.0
11	0	1	1	1.84	1.84	0.0	1.84
12	0	1	1	5.12	5.12	0.0	5.12
13	1	1	2	2.54	2.72	0.18	2.54
14	1	1	2	2.06	3.04	0.98	2.06
15	1	0	1	1.03	1.03	0.0	1.03
16	0	1	1	2.47	2.47	0.0	2.47
17	0	0	0	0.0	0.0	0.0	0.0
18	1	1	2	1.75	1.84	0.09	1.75
19	1	0	1	1.21	1.21	0.0	1.21
20	0	1	1	1.48	1.48	0.0	1.48

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	
CHI2	CHI1	Total	Count ¹	%
1	0	1	1	5.0
0	0	0	2	10.0
0	0	0	3	15.0
0	0	0	4	20.0
0	0	0	5	25.0
0	0	0	6	30.0
1	0	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

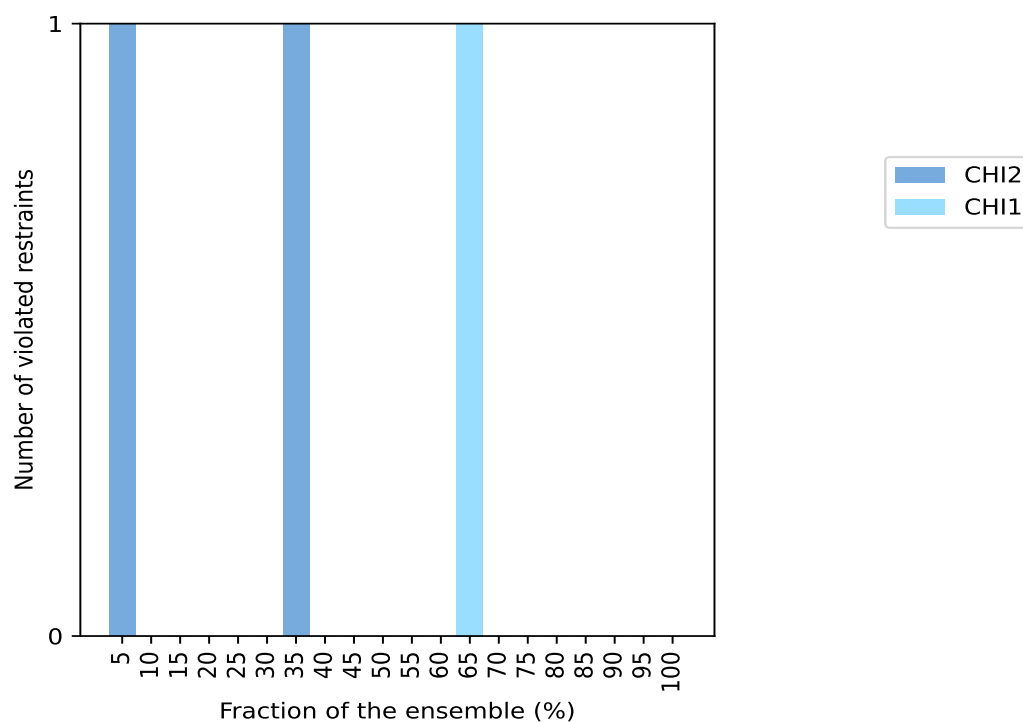
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Number of violated restraints			Fraction of the ensemble	
CHI2	CHI1	Total	Count ¹	%
0	0	0	12	60.0
0	1	1	13	65.0
0	0	0	14	70.0
0	0	0	15	75.0
0	0	0	16	80.0
0	0	0	17	85.0
0	0	0	18	90.0
0	0	0	19	95.0
0	0	0	20	100.0

¹ Number of models with violations

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

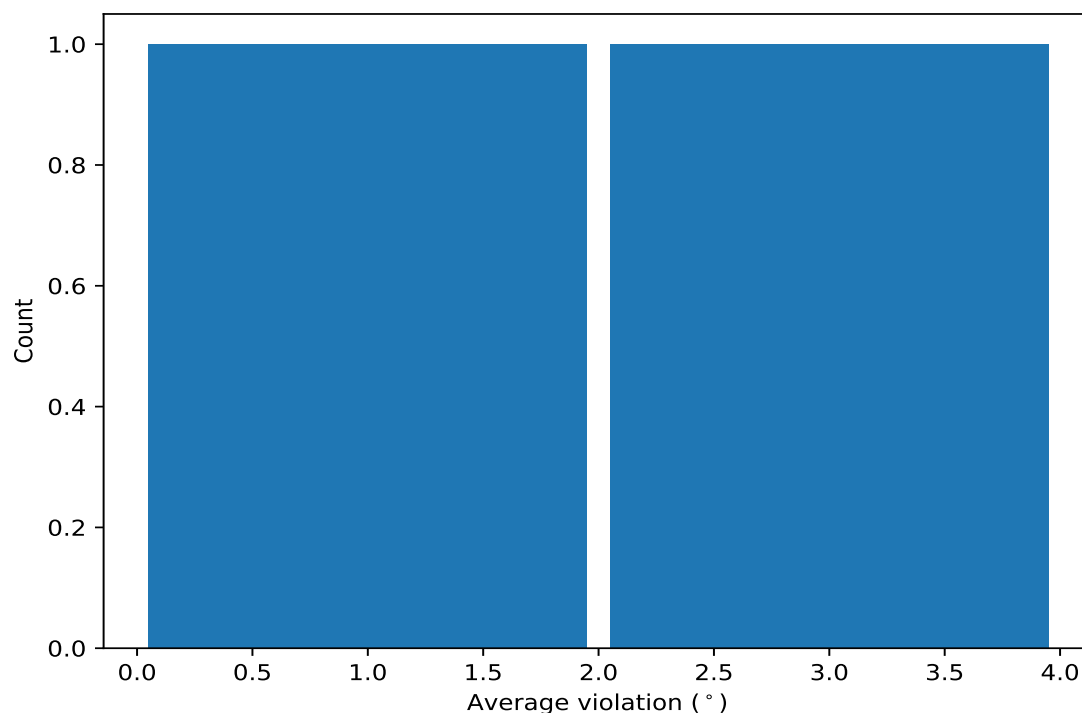


10.4 Most violated dihedral-angle restraints in the ensemble [i](#)

10.4.1 Histogram : Distribution of mean dihedral-angle violations [i](#)

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models

in the ensemble



10.4.2 Table: Most violated dihedral-angle restraints [i](#)

The following table provides the mean and the standard deviation of the violation for each restraint sorted by number of violated models and the mean value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint.

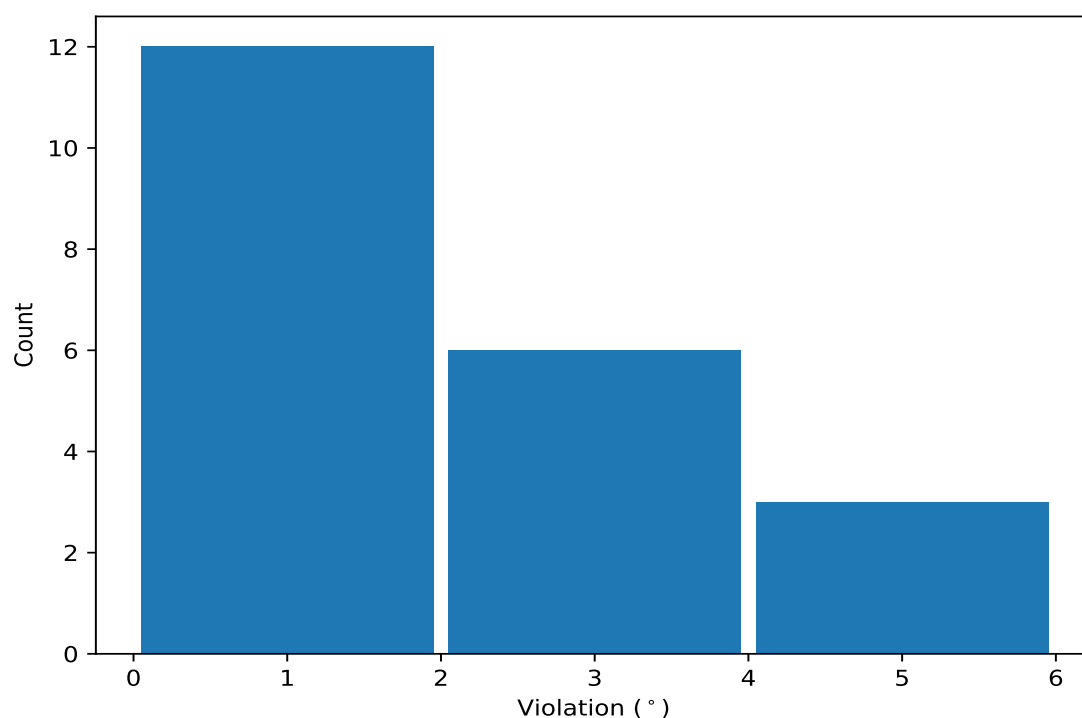
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models ¹	Mean	SD ²	Median
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	13	2.87	1.33	2.53
(1,10)	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1:41:A:HIS:ND1	7	1.59	0.46	1.74

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 (°)
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	9	5.4
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	12	5.12
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	3	4.21
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	8	3.91
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	14	3.04
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	13	2.72
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	7	2.53
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	16	2.47
(1,10)	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1:41:A:HIS:ND1	13	2.36
(1,10)	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1:41:A:HIS:ND1	5	1.86
(1,10)	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1:41:A:HIS:ND1	18	1.84
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	11	1.84
(1,10)	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1:41:A:HIS:ND1	6	1.74
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	18	1.66
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1	1.58
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	20	1.48
(1,9)	1:41:A:HIS:N	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	6	1.4
(1,8)	1:39:A:HIS:CA	1:39:A:HIS:CB	1:39:A:HIS:CG	1:39:A:HIS:ND1	2	1.25
(1,10)	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1:41:A:HIS:ND1	19	1.21
(1,10)	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1:41:A:HIS:ND1	14	1.07
(1,10)	1:41:A:HIS:CA	1:41:A:HIS:CB	1:41:A:HIS:CG	1:41:A:HIS:ND1	15	1.03