



# Full wwPDB NMR Structure Validation Report ⓘ

Jun 5, 2023 – 06:11 PM JST

PDB ID : 7BQQ  
BMRB ID : 36337  
Title : Solution NMR structure of fold-Z Gogy; de novo designed protein with an asymmetric all-alpha topology  
Authors : Kobayashi, N.; Sugiki, T.; Fujiwara, T.; Sakuma, K.; Kosugi, T.; Koga, R.; Koga, N.  
Deposited on : 2020-03-25

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
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) : 2.33

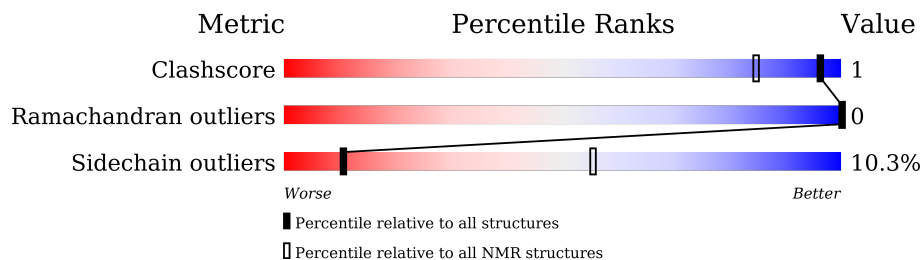
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 89%.

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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  $\geq 3$ , 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	130	

## 2 Ensemble composition and analysis i

This entry contains 20 models. Model 10 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *target function*.

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:9-A:115 (107)	0.29	10

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

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

The models can be grouped into 4 clusters and 3 single-model clusters were found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2129 atoms, of which 1061 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Gogy.

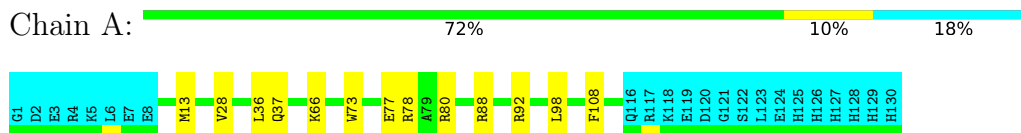
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	130	2129	648	1061	205	209	6	0

## 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: Gogy

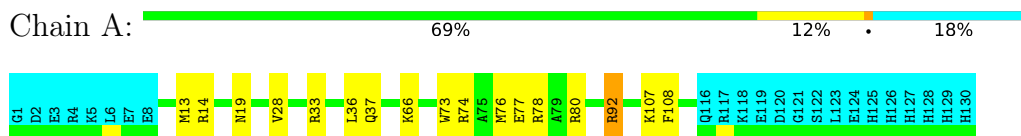


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

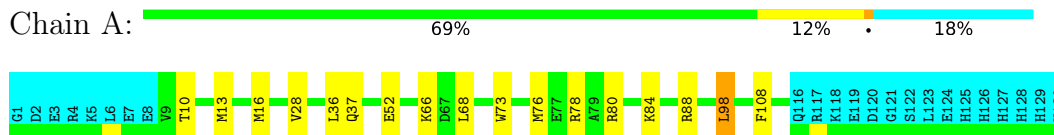
#### 4.2.1 Score per residue for model 1

- Molecule 1: Gogy



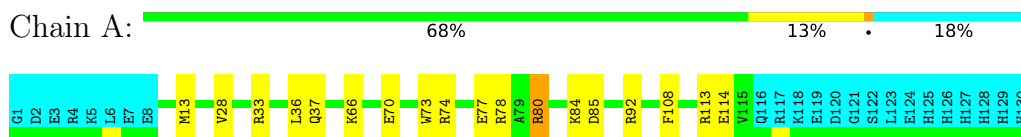
#### 4.2.2 Score per residue for model 2

- Molecule 1: Gogy



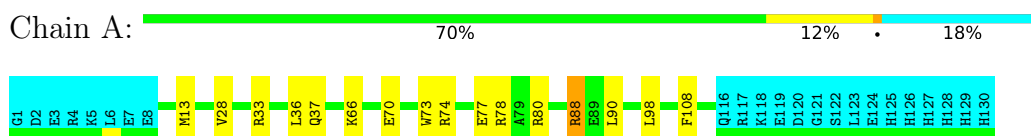
### 4.2.3 Score per residue for model 3

- Molecule 1: Gogy



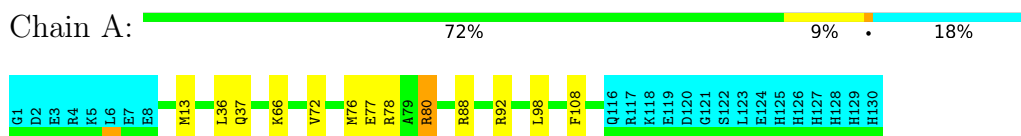
### 4.2.4 Score per residue for model 4

- Molecule 1: Gogy



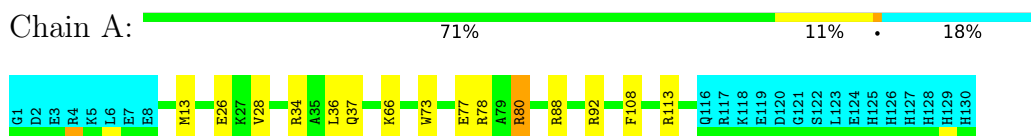
### 4.2.5 Score per residue for model 5

- Molecule 1: Gogy



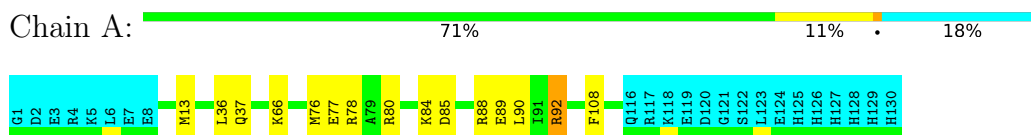
### 4.2.6 Score per residue for model 6

- Molecule 1: Gogy



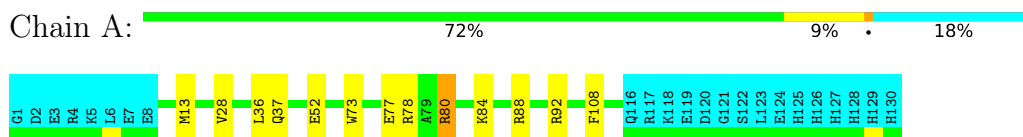
### 4.2.7 Score per residue for model 7

- Molecule 1: Gogy



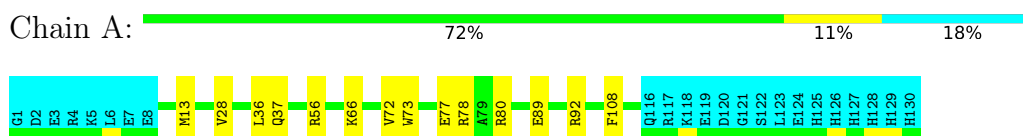
### 4.2.8 Score per residue for model 8

- Molecule 1: Gogy



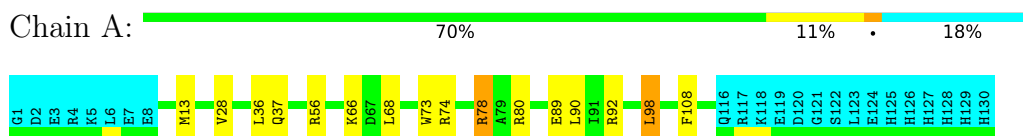
### 4.2.9 Score per residue for model 9

- Molecule 1: Gogy



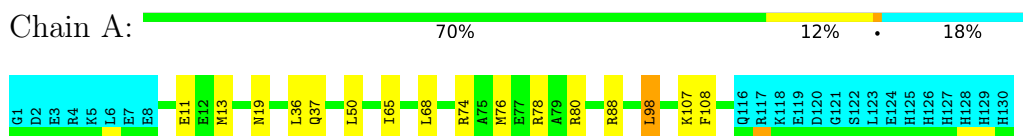
### 4.2.10 Score per residue for model 10 (medoid)

- Molecule 1: Gogy



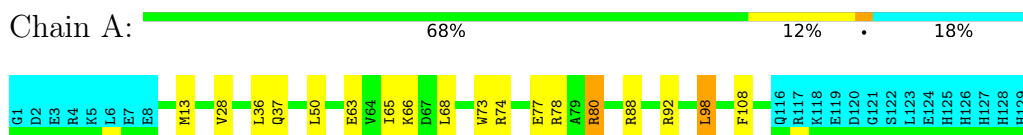
### 4.2.11 Score per residue for model 11

- Molecule 1: Gogy



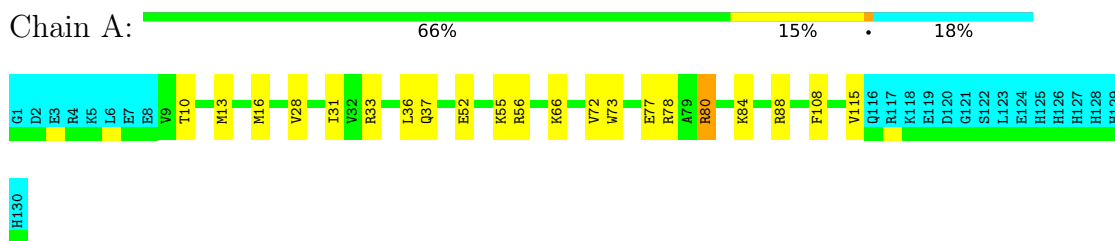
### 4.2.12 Score per residue for model 12

- Molecule 1: Gogy



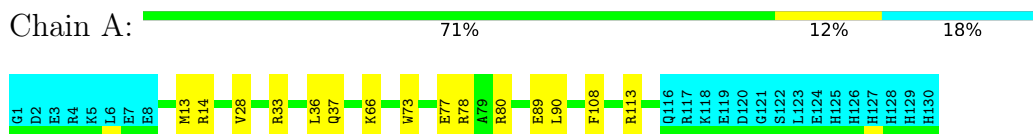
### 4.2.13 Score per residue for model 13

- Molecule 1: Gogy



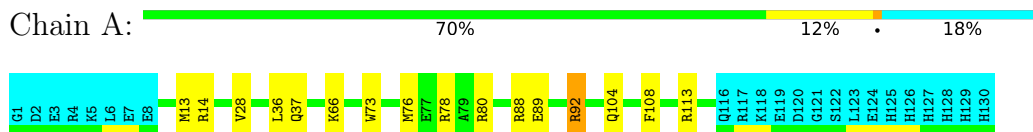
### 4.2.14 Score per residue for model 14

- Molecule 1: Gogy



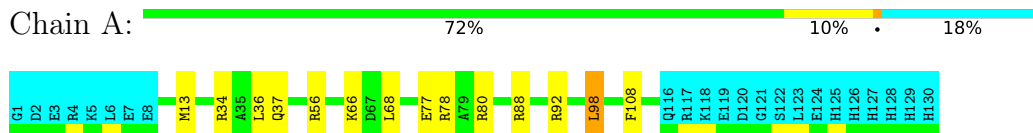
### 4.2.15 Score per residue for model 15

- Molecule 1: Gogy



### 4.2.16 Score per residue for model 16

- Molecule 1: Gogy

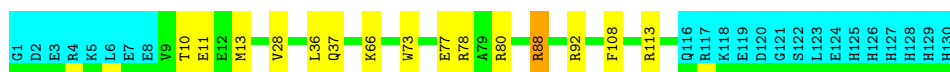


### 4.2.17 Score per residue for model 17

- Molecule 1: Gogy







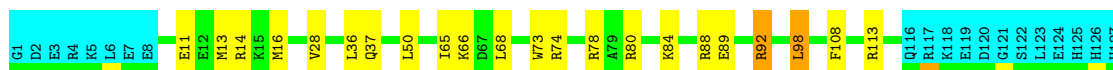
#### 4.2.18 Score per residue for model 18

- Molecule 1: Gogy



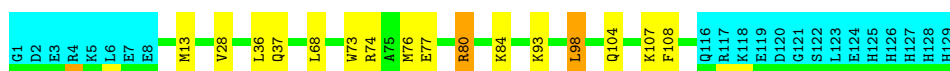
#### 4.2.19 Score per residue for model 19

- Molecule 1: Gogy



#### 4.2.20 Score per residue for model 20

- Molecule 1: Gogy



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *molecular dynamics*.

Of the 200 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 calculation	3.98
Amber	refinement	12

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	1556
Number of shifts mapped to atoms	1556
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	89%

## 6 Model quality i

### 6.1 Standard geometry i

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	0.65±0.00	0±0/875 ( 0.0± 0.0%)	1.06±0.03	5±1/1168 ( 0.4± 0.1%)
All	All	0.65	0/17500 ( 0.0%)	1.06	102/23360 ( 0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	0.2±0.4
All	All	0	4

There are no bond-length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	88	ARG	NE-CZ-NH1	10.22	125.41	120.30	7	14
1	A	80	ARG	NE-CZ-NH1	8.54	124.57	120.30	4	20
1	A	78	ARG	NE-CZ-NH1	8.46	124.53	120.30	6	19
1	A	74	ARG	NE-CZ-NH1	6.98	123.79	120.30	20	7
1	A	92	ARG	NE-CZ-NH1	6.94	123.77	120.30	3	13
1	A	113	ARG	NE-CZ-NH1	6.93	123.77	120.30	14	5
1	A	33	ARG	NE-CZ-NH1	6.87	123.74	120.30	4	5
1	A	14	ARG	NE-CZ-NH1	6.58	123.59	120.30	15	4
1	A	56	ARG	NE-CZ-NH1	6.30	123.45	120.30	10	4
1	A	74	ARG	NE-CZ-NH2	-6.21	117.19	120.30	4	3
1	A	34	ARG	NE-CZ-NH1	6.12	123.36	120.30	16	2
1	A	88	ARG	NE-CZ-NH2	-5.93	117.33	120.30	7	1
1	A	88	ARG	CD-NE-CZ	5.77	131.68	123.60	7	1
1	A	78	ARG	NE-CZ-NH2	-5.32	117.64	120.30	9	2
1	A	92	ARG	NE-CZ-NH2	-5.24	117.68	120.30	5	2

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	92	ARG	Sidechain	2
1	A	74	ARG	Sidechain	2

## 6.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	867	886	886	1±1
All	All	17340	17720	17720	29

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 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:28:VAL:HG11	1:A:73:TRP:CE3	0.54	2.37	3	16
1:A:68:LEU:HD11	1:A:98:LEU:HD13	0.51	1.83	19	5
1:A:50:LEU:HD13	1:A:65:ILE:HD11	0.47	1.86	12	3
1:A:31:ILE:HG23	1:A:115:VAL:HG23	0.43	1.90	13	1
1:A:68:LEU:HD11	1:A:98:LEU:CD1	0.42	2.44	10	4

## 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	107/130 (82%)	106±1 (99±1%)	1±1 (1±1%)	0±0 (0±0%)	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2140/2600 (82%)	2124 (99%)	16 (1%)	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	94/115 (82%)	84±2 (90±2%)	10±2 (10±2%)	11	55
All	All	1880/2300 (82%)	1687 (90%)	193 (10%)	11	55

All 32 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	13	MET	20
1	A	36	LEU	20
1	A	37	GLN	20
1	A	108	PHE	20
1	A	66	LYS	17
1	A	77	GLU	15
1	A	98	LEU	9
1	A	76	MET	8
1	A	84	LYS	7
1	A	80	ARG	7
1	A	89	GLU	7
1	A	90	LEU	4
1	A	92	ARG	3
1	A	107	LYS	3
1	A	10	THR	3
1	A	16	MET	3
1	A	52	GLU	3
1	A	72	VAL	3
1	A	11	GLU	3
1	A	19	ASN	2
1	A	70	GLU	2
1	A	85	ASP	2

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Mol	Chain	Res	Type	Models (Total)
1	A	88	ARG	2
1	A	104	GLN	2
1	A	33	ARG	1
1	A	113	ARG	1
1	A	114	GLU	1
1	A	26	GLU	1
1	A	78	ARG	1
1	A	63	GLU	1
1	A	55	LYS	1
1	A	93	LYS	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 monosaccharides in this entry.

### 6.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 6.7 Other polymers [i](#)

There are no such molecules in this entry.

### 6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 7 Chemical shift validation [i](#)

The completeness of assignment taking into account all chemical shift lists is 89% for the well-defined parts and 84% for the entire structure.

### 7.1 Chemical shift list 1

File name: working\_cs.cif

Chemical shift list name: *shift\_set\_1*

#### 7.1.1 Bookkeeping [i](#)

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

Total number of shifts	1556
Number of shifts mapped to atoms	1556
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

#### 7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	124	$-0.49 \pm 0.10$	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	117	$0.19 \pm 0.06$	None needed (< 0.5 ppm)
$^{13}\text{C}'$	122	$-0.37 \pm 0.06$	None needed (< 0.5 ppm)
$^{15}\text{N}$	118	$-0.17 \pm 0.31$	None needed (< 0.5 ppm)

#### 7.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 89%, i.e. 1364 atoms were assigned a chemical shift out of a possible 1529. 0 out of 14 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	531/531 (100%)	214/214 (100%)	214/214 (100%)	103/103 (100%)
Sidechain	801/966 (83%)	550/616 (89%)	242/297 (81%)	9/53 (17%)

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	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Aromatic	32/32 (100%)	16/16 (100%)	15/15 (100%)	1/1 (100%)
Overall	1364/1529 (89%)	780/846 (92%)	471/526 (90%)	113/157 (72%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 84%, i.e. 1554 atoms were assigned a chemical shift out of a possible 1850. 0 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	611/648 (94%)	247/262 (94%)	246/260 (95%)	118/126 (94%)
Sidechain	911/1128 (81%)	624/716 (87%)	277/350 (79%)	10/62 (16%)
Aromatic	32/74 (43%)	16/40 (40%)	15/27 (56%)	1/7 (14%)
Overall	1554/1850 (84%)	887/1018 (87%)	538/637 (84%)	129/195 (66%)

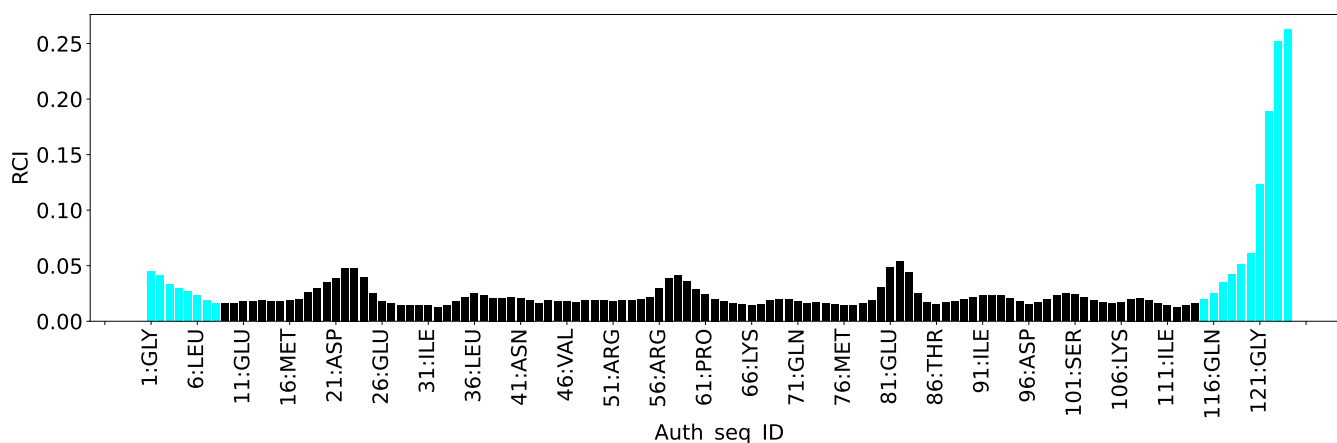
#### 7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

#### 7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:





## 8 NMR restraints analysis

### 8.1 Conformationally restricting restraints

The following table provides the summary of experimentally observed NMR restraints in different categories. Restraints are classified into different categories based on the sequence separation of the atoms involved.

Description	Value
Total distance restraints	2639
Intra-residue ( $ i-j =0$ )	509
Sequential ( $ i-j =1$ )	593
Medium range ( $ i-j >1$ and $ i-j <5$ )	841
Long range ( $ i-j \geq 5$ )	696
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	20.3
Number of long range restraints per residue <sup>1</sup>	5.4

<sup>1</sup>Long 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)	8.0	0.2
0.2-0.5 (Medium)	0.4	0.31
>0.5 (Large)	None	None

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

Dihedral-angle violations less than  $1^\circ$  are not included in the calculation. There are no dihedral-angle violations

## 9 Distance violation analysis [i](#)

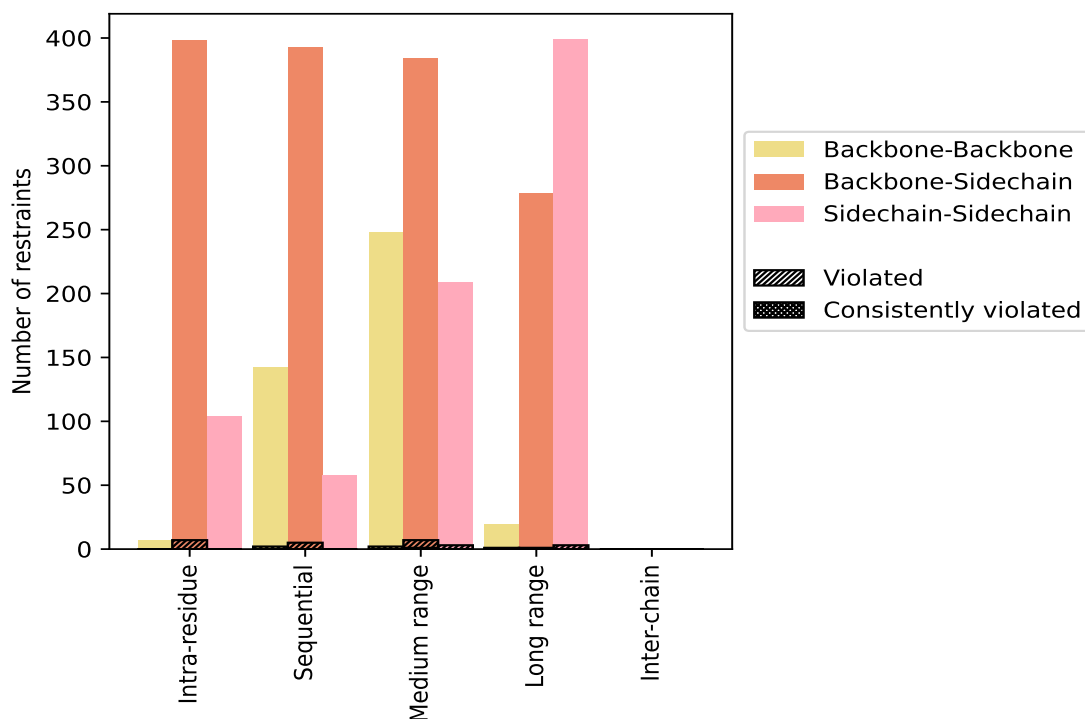
### 9.1 Summary of distance violations [i](#)

The following table shows the summary of distance violations in different restraint categories based on the sequence separation of the atoms involved. Each category is further sub-divided into three sub-categories based on the atoms involved. Violations less than 0.1 Å are not included in the statistics.

Restrains type	Count	% <sup>1</sup>	Violated <sup>3</sup>			Consistently Violated <sup>4</sup>		
			Count	% <sup>2</sup>	% <sup>1</sup>	Count	% <sup>2</sup>	% <sup>1</sup>
<b>Intra-residue ( i-j =0)</b>	<b>509</b>	<b>19.3</b>	<b>7</b>	<b>1.4</b>	<b>0.3</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	7	0.3	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	398	15.1	7	1.8	0.3	0	0.0	0.0
Sidechain-Sidechain	104	3.9	0	0.0	0.0	0	0.0	0.0
<b>Sequential ( i-j =1)</b>	<b>593</b>	<b>22.5</b>	<b>7</b>	<b>1.2</b>	<b>0.3</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	142	5.4	2	1.4	0.1	0	0.0	0.0
Backbone-Sidechain	393	14.9	5	1.3	0.2	0	0.0	0.0
Sidechain-Sidechain	58	2.2	0	0.0	0.0	0	0.0	0.0
<b>Medium range ( i-j &gt;1 &amp;  i-j &lt;5)</b>	<b>841</b>	<b>31.9</b>	<b>12</b>	<b>1.4</b>	<b>0.5</b>	<b>1</b>	<b>0.1</b>	<b>0.0</b>
Backbone-Backbone	248	9.4	2	0.8	0.1	0	0.0	0.0
Backbone-Sidechain	384	14.6	7	1.8	0.3	1	0.3	0.0
Sidechain-Sidechain	209	7.9	3	1.4	0.1	0	0.0	0.0
<b>Long range ( i-j ≥5)</b>	<b>696</b>	<b>26.4</b>	<b>5</b>	<b>0.7</b>	<b>0.2</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	19	0.7	1	5.3	0.0	0	0.0	0.0
Backbone-Sidechain	278	10.5	1	0.4	0.0	0	0.0	0.0
Sidechain-Sidechain	399	15.1	3	0.8	0.1	0	0.0	0.0
<b>Inter-chain</b>	<b>0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
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
<b>Hydrogen bond</b>	<b>0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
<b>Disulfide bond</b>	<b>0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
<b>Total</b>	<b>2639</b>	<b>100.0</b>	<b>31</b>	<b>1.2</b>	<b>1.2</b>	<b>1</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	416	15.8	5	1.2	0.2	0	0.0	0.0
Backbone-Sidechain	1453	55.1	20	1.4	0.8	1	0.1	0.0
Sidechain-Sidechain	770	29.2	6	0.8	0.2	0	0.0	0.0

<sup>1</sup> percentage calculated with respect to the total number of distance restraints, <sup>2</sup> percentage calculated with respect to the number of restraints in a particular restraint category, <sup>3</sup> violated in at least one model, <sup>4</sup> violated in all the models

### 9.1.1 Bar chart : Distribution of distance restraints and violations [i](#)



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and disulfid bonds are counted in their appropriate category on the x-axis

## 9.2 Distance violation statistics for each model [i](#)

The following table provides the distance violation statistics for each model in the ensemble. Violations less than 0.1 Å are not included in the statistics.

Model ID	Number of violations						Mean (Å)	Max (Å)	SD <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total				
1	1	1	3	1	0	6	0.14	0.17	0.02	0.14
2	2	1	6	1	0	10	0.13	0.2	0.03	0.12
3	1	0	5	1	0	7	0.13	0.17	0.02	0.13
4	2	1	4	3	0	10	0.13	0.2	0.03	0.12
5	0	0	5	1	0	6	0.14	0.2	0.03	0.14
6	1	2	5	2	0	10	0.13	0.15	0.01	0.12
7	1	0	4	2	0	7	0.12	0.18	0.02	0.12
8	1	1	3	2	0	7	0.13	0.17	0.02	0.12
9	2	0	3	2	0	7	0.12	0.14	0.01	0.12
10	2	1	6	2	0	11	0.14	0.2	0.03	0.12
11	1	1	5	2	0	9	0.14	0.21	0.03	0.13

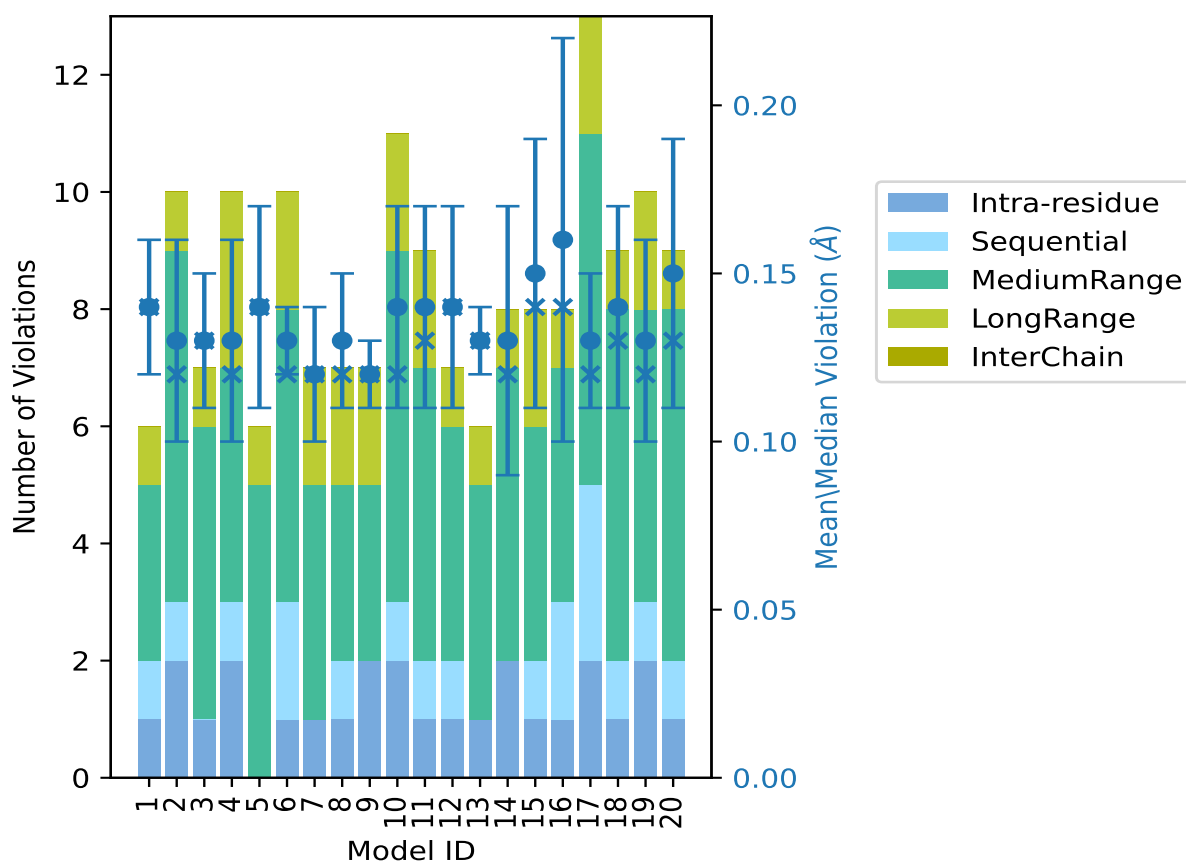
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Model ID	Number of violations					Total	Mean (Å)	Max (Å)	SD <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>					
12	1	1	4	1	0	7	0.14	0.21	0.03	0.14
13	1	0	4	1	0	6	0.13	0.15	0.01	0.13
14	2	0	5	1	0	8	0.13	0.23	0.04	0.12
15	1	1	4	2	0	8	0.15	0.22	0.04	0.14
16	1	2	4	1	0	8	0.16	0.31	0.06	0.14
17	2	3	6	2	0	13	0.13	0.17	0.02	0.12
18	1	1	6	1	0	9	0.14	0.19	0.03	0.13
19	2	1	5	2	0	10	0.13	0.22	0.03	0.12
20	1	1	6	1	0	9	0.15	0.21	0.04	0.13

<sup>1</sup>Intra-residue restraints, <sup>2</sup>Sequential restraints, <sup>3</sup>Medium range restraints, <sup>4</sup>Long range restraints, <sup>5</sup>Inter-chain restraints, <sup>6</sup>Standard deviation

### 9.2.1 Bar graph : Distance Violation statistics for each model [i](#)



The mean(dot), median(x) and the standard deviation are shown in blue with respect to the y axis on the right

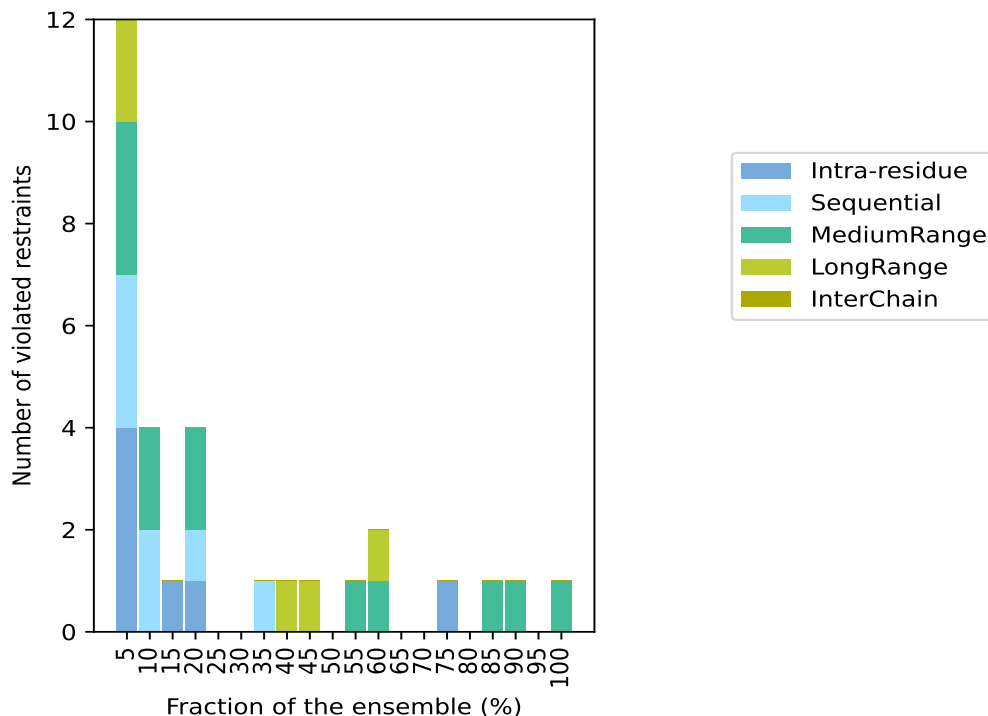
### 9.3 Distance violation statistics for the ensemble [i](#)

Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 2608(IR:502, SQ:586, MR:829, LR:691, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total	Count <sup>6</sup>	%
4	3	3	2	0	12	1	5.0
0	2	2	0	0	4	2	10.0
1	0	0	0	0	1	3	15.0
1	1	2	0	0	4	4	20.0
0	0	0	0	0	0	5	25.0
0	0	0	0	0	0	6	30.0
0	1	0	0	0	1	7	35.0
0	0	0	1	0	1	8	40.0
0	0	0	1	0	1	9	45.0
0	0	0	0	0	0	10	50.0
0	0	1	0	0	1	11	55.0
0	0	1	1	0	2	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	1	0	0	1	17	85.0
0	0	1	0	0	1	18	90.0
0	0	0	0	0	0	19	95.0
0	0	1	0	0	1	20	100.0

<sup>1</sup>Intra-residue restraints, <sup>2</sup>Sequential restraints, <sup>3</sup>Medium range restraints, <sup>4</sup>Long range restraints, <sup>5</sup>Inter-chain restraints, <sup>6</sup> Number of models with violations

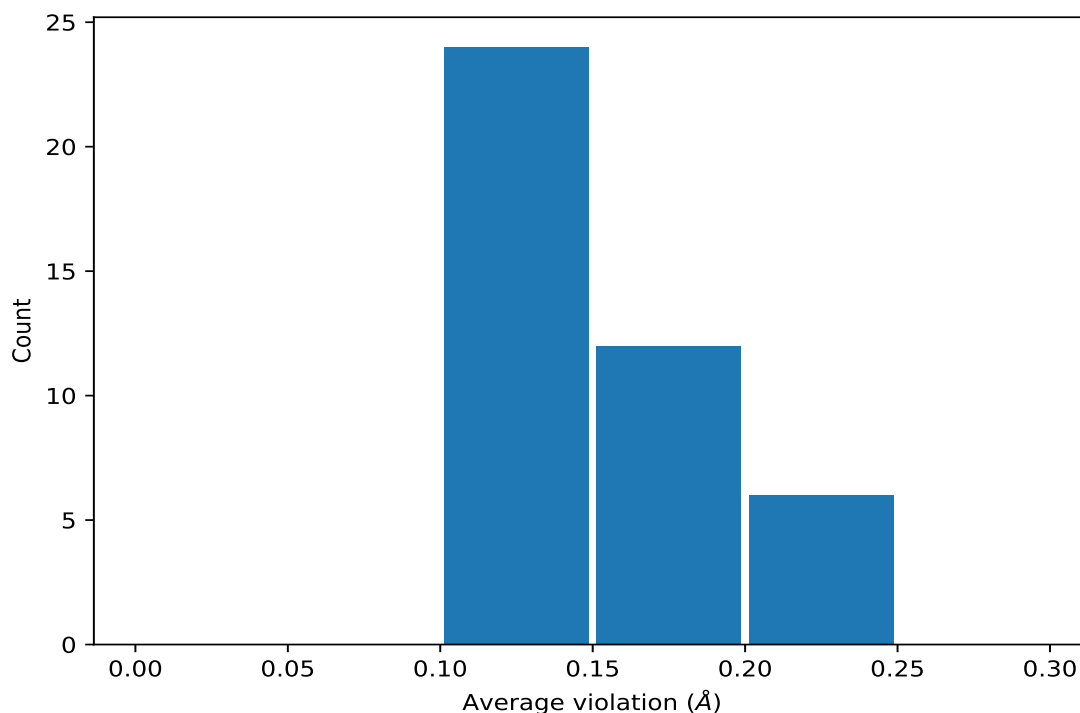
### 9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



## 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 <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	20	0.12	0.01	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	20	0.12	0.01	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	20	0.12	0.01	0.12
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	18	0.15	0.03	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	18	0.15	0.03	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	18	0.15	0.03	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	18	0.15	0.03	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	18	0.15	0.03	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	18	0.15	0.03	0.15
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	17	0.13	0.01	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	17	0.13	0.01	0.13
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	15	0.12	0.01	0.12
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	12	0.13	0.01	0.14
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	12	0.12	0.01	0.11
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	11	0.12	0.01	0.12
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	11	0.12	0.01	0.12

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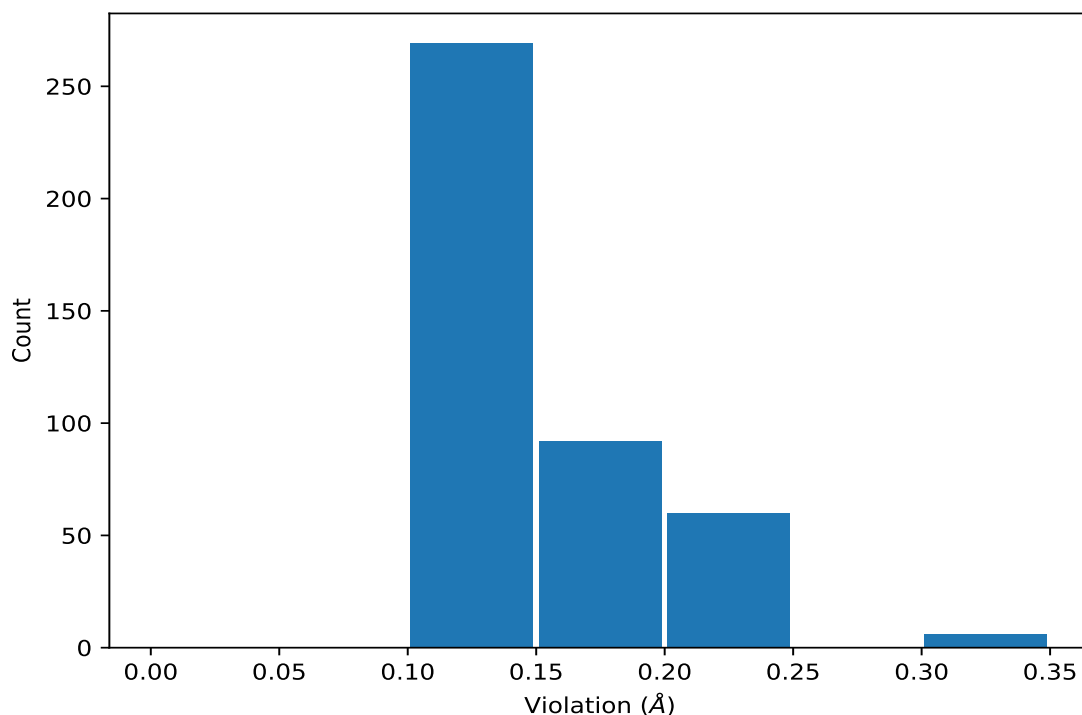
Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	9	0.21	0.01	0.21
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	9	0.21	0.01	0.21
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	9	0.21	0.01	0.21
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	9	0.21	0.01	0.21
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	9	0.21	0.01	0.21
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	9	0.21	0.01	0.21
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	8	0.14	0.02	0.15
(1,663)	1:A:38:GLN:HG3	1:A:39:MET:H	7	0.12	0.01	0.12
(1,1248)	1:A:3:GLU:HG2	1:A:4:ARG:H	4	0.16	0.01	0.16
(1,1248)	1:A:3:GLU:HG3	1:A:4:ARG:H	4	0.16	0.01	0.16
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE1	4	0.12	0.01	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE2	4	0.12	0.01	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE3	4	0.12	0.01	0.11
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE2	4	0.11	0.0	0.11
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE3	4	0.11	0.0	0.11
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE2	4	0.11	0.0	0.11
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE3	4	0.11	0.0	0.11
(1,926)	1:A:79:ALA:H	1:A:81:GLU:H	4	0.11	0.0	0.11
(1,2427)	1:A:56:ARG:H	1:A:56:ARG:HG2	3	0.13	0.01	0.13
(1,2427)	1:A:56:ARG:H	1:A:56:ARG:HG3	3	0.13	0.01	0.13
(1,1365)	1:A:115:VAL:HB	1:A:118:LYS:HB2	2	0.19	0.01	0.19
(1,1365)	1:A:115:VAL:HB	1:A:118:LYS:HB3	2	0.19	0.01	0.19
(1,2638)	1:A:123:LEU:HB2	1:A:124:GLU:H	2	0.19	0.03	0.19
(1,2638)	1:A:123:LEU:HB3	1:A:124:GLU:H	2	0.19	0.03	0.19
(1,66)	1:A:2:ASP:HA	1:A:3:GLU:H	2	0.14	0.03	0.14
(1,614)	1:A:68:LEU:H	1:A:71:GLN:H	2	0.11	0.0	0.11

<sup>1</sup>Number of violated models, <sup>2</sup>Standard 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)
(1,2637)	1:A:123:LEU:HA	1:A:123:LEU:HD11	16	0.31
(1,2637)	1:A:123:LEU:HA	1:A:123:LEU:HD12	16	0.31
(1,2637)	1:A:123:LEU:HA	1:A:123:LEU:HD13	16	0.31
(1,2637)	1:A:123:LEU:HA	1:A:123:LEU:HD21	16	0.31
(1,2637)	1:A:123:LEU:HA	1:A:123:LEU:HD22	16	0.31
(1,2637)	1:A:123:LEU:HA	1:A:123:LEU:HD23	16	0.31
(1,2230)	1:A:11:GLU:H	1:A:11:GLU:HG2	14	0.23
(1,2230)	1:A:11:GLU:H	1:A:11:GLU:HG3	14	0.23
(1,2638)	1:A:123:LEU:HB2	1:A:124:GLU:H	15	0.22
(1,2638)	1:A:123:LEU:HB3	1:A:124:GLU:H	15	0.22
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	16	0.22
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	16	0.22
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	16	0.22
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	16	0.22
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	16	0.22
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	16	0.22

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	19	0.22
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	19	0.22
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	19	0.22
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	19	0.22
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	19	0.22
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	19	0.22
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	11	0.21
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	11	0.21
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	11	0.21
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	11	0.21
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	11	0.21
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	11	0.21
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	12	0.21
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	12	0.21
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	12	0.21
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	12	0.21
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	12	0.21
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	12	0.21
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	20	0.21
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	20	0.21
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	20	0.21
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	20	0.21
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	20	0.21
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	20	0.21
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	2	0.2
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	2	0.2
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	2	0.2
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	2	0.2
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	2	0.2
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	2	0.2
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	4	0.2
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	4	0.2
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	4	0.2
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	4	0.2
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	4	0.2
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	4	0.2
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	5	0.2
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	5	0.2
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	5	0.2
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	5	0.2
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	5	0.2
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	5	0.2

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2463)	1:A:68:LEU:HD11	1:A:98:LEU:HG	10	0.2
(1,2463)	1:A:68:LEU:HD12	1:A:98:LEU:HG	10	0.2
(1,2463)	1:A:68:LEU:HD13	1:A:98:LEU:HG	10	0.2
(1,2463)	1:A:68:LEU:HD21	1:A:98:LEU:HG	10	0.2
(1,2463)	1:A:68:LEU:HD22	1:A:98:LEU:HG	10	0.2
(1,2463)	1:A:68:LEU:HD23	1:A:98:LEU:HG	10	0.2
(1,1365)	1:A:115:VAL:HB	1:A:118:LYS:HB2	20	0.2
(1,1365)	1:A:115:VAL:HB	1:A:118:LYS:HB3	20	0.2
(1,473)	1:A:50:LEU:H	1:A:51:ARG:HG2	18	0.19
(1,473)	1:A:50:LEU:H	1:A:51:ARG:HG3	18	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	15	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	15	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	15	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	15	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	15	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	15	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	20	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	20	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	20	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	20	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	20	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	20	0.19
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	7	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	7	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	7	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	7	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	7	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	7	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	18	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	18	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	18	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	18	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	18	0.18
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	18	0.18
(1,1365)	1:A:115:VAL:HB	1:A:118:LYS:HB2	10	0.18
(1,1365)	1:A:115:VAL:HB	1:A:118:LYS:HB3	10	0.18
(1,66)	1:A:2:ASP:HA	1:A:3:GLU:H	17	0.17
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	8	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	2	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	2	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	2	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	2	0.17

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	2	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	2	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	3	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	3	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	3	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	3	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	3	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	3	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	19	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	19	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	19	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	19	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	19	0.17
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	19	0.17
(1,1248)	1:A:3:GLU:HG2	1:A:4:ARG:H	1	0.17
(1,1248)	1:A:3:GLU:HG3	1:A:4:ARG:H	1	0.17
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	15	0.16
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	17	0.16
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	17	0.16
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	17	0.16
(1,2638)	1:A:123:LEU:HB2	1:A:124:GLU:H	10	0.16
(1,2638)	1:A:123:LEU:HB3	1:A:124:GLU:H	10	0.16
(1,1248)	1:A:3:GLU:HG2	1:A:4:ARG:H	16	0.16
(1,1248)	1:A:3:GLU:HG3	1:A:4:ARG:H	16	0.16
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	18	0.15
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	18	0.15
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	6	0.15
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	11	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	12	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	12	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	12	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	12	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	12	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	12	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	13	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	13	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	13	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	13	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	13	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	13	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	16	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	16	0.15

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	16	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	16	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	16	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	16	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	17	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	17	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	17	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	17	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	17	0.15
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	17	0.15
(1,1258)	1:A:3:GLU:HA	1:A:3:GLU:HG2	17	0.15
(1,1258)	1:A:3:GLU:HA	1:A:3:GLU:HG3	17	0.15
(1,1248)	1:A:3:GLU:HG2	1:A:4:ARG:H	4	0.15
(1,1248)	1:A:3:GLU:HG3	1:A:4:ARG:H	4	0.15
(1,1248)	1:A:3:GLU:HG2	1:A:4:ARG:H	8	0.15
(1,1248)	1:A:3:GLU:HG3	1:A:4:ARG:H	8	0.15
(1,663)	1:A:38:GLN:HG3	1:A:39:MET:H	12	0.14
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	1	0.14
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	1	0.14
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	5	0.14
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	5	0.14
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	20	0.14
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	20	0.14
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	9	0.14
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	5	0.14
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	5	0.14
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	5	0.14
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	12	0.14
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	12	0.14
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	12	0.14
(1,2596)	1:A:112:MET:H	1:A:114:GLU:HB2	3	0.14
(1,2596)	1:A:112:MET:H	1:A:114:GLU:HB3	3	0.14
(1,2427)	1:A:56:ARG:H	1:A:56:ARG:HG2	15	0.14
(1,2427)	1:A:56:ARG:H	1:A:56:ARG:HG3	15	0.14
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	1	0.14
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	1	0.14
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	1	0.14
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	1	0.14
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	1	0.14
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	1	0.14
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	1	0.14
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	3	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	6	0.14
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	13	0.14
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	14	0.14
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	17	0.14
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	18	0.14
(1,663)	1:A:38:GLN:HG3	1:A:39:MET:H	11	0.13
(1,663)	1:A:38:GLN:HG3	1:A:39:MET:H	20	0.13
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	2	0.13
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	2	0.13
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	3	0.13
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	3	0.13
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	16	0.13
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	16	0.13
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	17	0.13
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	17	0.13
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	20	0.13
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	20	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	2	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	2	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	4	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	4	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	6	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	6	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	9	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	9	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	13	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	13	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	15	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	15	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	16	0.13
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	16	0.13
(1,385)	1:A:123:LEU:HA	1:A:124:GLU:H	6	0.13
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	6	0.13
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	6	0.13
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	6	0.13
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	11	0.13
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	11	0.13
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	11	0.13
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	13	0.13
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	13	0.13
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	13	0.13
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	15	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	15	0.13
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	15	0.13
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	20	0.13
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	20	0.13
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	20	0.13
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	2	0.13
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	11	0.13
(1,2427)	1:A:56:ARG:H	1:A:56:ARG:HG2	19	0.13
(1,2427)	1:A:56:ARG:H	1:A:56:ARG:HG3	19	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	4	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	4	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	4	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	4	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	4	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	4	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	5	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	5	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	5	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	5	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	5	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	5	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	10	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	10	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	10	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	10	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	10	0.13
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	10	0.13
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE1	18	0.13
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE2	18	0.13
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE3	18	0.13
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	11	0.13
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	12	0.13
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	8	0.13
(1,671)	1:A:3:GLU:HB2	1:A:5:LYS:H	17	0.12
(1,663)	1:A:38:GLN:HG3	1:A:39:MET:H	2	0.12
(1,663)	1:A:38:GLN:HG3	1:A:39:MET:H	6	0.12
(1,663)	1:A:38:GLN:HG3	1:A:39:MET:H	17	0.12
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	6	0.12
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	6	0.12
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	9	0.12
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	9	0.12
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	10	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	10	0.12
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	19	0.12
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	19	0.12
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	10	0.12
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	10	0.12
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	11	0.12
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	11	0.12
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	14	0.12
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	14	0.12
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	17	0.12
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	17	0.12
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	4	0.12
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	7	0.12
(1,544)	1:A:12:GLU:H	1:A:106:LYS:HA	10	0.12
(1,410)	1:A:12:GLU:HB2	1:A:13:MET:H	17	0.12
(1,410)	1:A:12:GLU:HB3	1:A:13:MET:H	17	0.12
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	2	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	2	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	2	0.12
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	3	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	3	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	3	0.12
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	8	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	8	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	8	0.12
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	9	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	9	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	9	0.12
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	10	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	10	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	10	0.12
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	14	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	14	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	14	0.12
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	18	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	18	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	18	0.12
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	19	0.12
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	19	0.12
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	19	0.12
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	1	0.12
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	3	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	4	0.12
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	7	0.12
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	9	0.12
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	10	0.12
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	13	0.12
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	17	0.12
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	19	0.12
(1,2427)	1:A:56:ARG:H	1:A:56:ARG:HG2	12	0.12
(1,2427)	1:A:56:ARG:H	1:A:56:ARG:HG3	12	0.12
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	14	0.12
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	14	0.12
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	14	0.12
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	14	0.12
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	14	0.12
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	14	0.12
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	8	0.12
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	14	0.12
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	19	0.12
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	7	0.12
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	9	0.12
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	15	0.12
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE2	18	0.12
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE3	18	0.12
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE2	18	0.12
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE3	18	0.12
(1,926)	1:A:79:ALA:H	1:A:81:GLU:H	4	0.11
(1,926)	1:A:79:ALA:H	1:A:81:GLU:H	7	0.11
(1,926)	1:A:79:ALA:H	1:A:81:GLU:H	10	0.11
(1,926)	1:A:79:ALA:H	1:A:81:GLU:H	14	0.11
(1,680)	1:A:10:THR:H	1:A:10:THR:HB	2	0.11
(1,663)	1:A:38:GLN:HG3	1:A:39:MET:H	19	0.11
(1,66)	1:A:2:ASP:HA	1:A:3:GLU:H	16	0.11
(1,614)	1:A:68:LEU:H	1:A:71:GLN:H	11	0.11
(1,614)	1:A:68:LEU:H	1:A:71:GLN:H	15	0.11
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	5	0.11
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	5	0.11
(1,584)	1:A:78:ARG:HD2	1:A:80:ARG:H	18	0.11
(1,584)	1:A:78:ARG:HD3	1:A:80:ARG:H	18	0.11
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	12	0.11
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	12	0.11
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG2	19	0.11
(1,58)	1:A:35:ALA:H	1:A:37:GLN:HG3	19	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	1	0.11
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	1	0.11
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	1	0.11
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	4	0.11
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	4	0.11
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	4	0.11
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	7	0.11
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	7	0.11
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	7	0.11
(1,311)	1:A:91:ILE:HD11	1:A:95:ALA:H	16	0.11
(1,311)	1:A:91:ILE:HD12	1:A:95:ALA:H	16	0.11
(1,311)	1:A:91:ILE:HD13	1:A:95:ALA:H	16	0.11
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	6	0.11
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	8	0.11
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	14	0.11
(1,265)	1:A:64:VAL:H	1:A:64:VAL:HB	20	0.11
(1,2197)	1:A:6:LEU:HD11	1:A:9:VAL:H	6	0.11
(1,2197)	1:A:6:LEU:HD12	1:A:9:VAL:H	6	0.11
(1,2197)	1:A:6:LEU:HD13	1:A:9:VAL:H	6	0.11
(1,2197)	1:A:6:LEU:HD21	1:A:9:VAL:H	6	0.11
(1,2197)	1:A:6:LEU:HD22	1:A:9:VAL:H	6	0.11
(1,2197)	1:A:6:LEU:HD23	1:A:9:VAL:H	6	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	8	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	8	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	8	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	8	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	8	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	8	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG11	11	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG12	11	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG13	11	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG21	11	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG22	11	0.11
(1,2195)	1:A:6:LEU:HG	1:A:9:VAL:HG23	11	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE1	4	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE2	4	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE3	4	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE1	9	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE2	9	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE3	9	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE1	10	0.11
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE2	10	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1987)	1:A:39:MET:H	1:A:39:MET:HE3	10	0.11
(1,1716)	1:A:32:VAL:HG11	1:A:71:GLN:HG2	17	0.11
(1,1716)	1:A:32:VAL:HG11	1:A:71:GLN:HG3	17	0.11
(1,1716)	1:A:32:VAL:HG12	1:A:71:GLN:HG2	17	0.11
(1,1716)	1:A:32:VAL:HG12	1:A:71:GLN:HG3	17	0.11
(1,1716)	1:A:32:VAL:HG13	1:A:71:GLN:HG2	17	0.11
(1,1716)	1:A:32:VAL:HG13	1:A:71:GLN:HG3	17	0.11
(1,1687)	1:A:6:LEU:HD21	1:A:88:ARG:HD2	4	0.11
(1,1687)	1:A:6:LEU:HD21	1:A:88:ARG:HD3	4	0.11
(1,1687)	1:A:6:LEU:HD22	1:A:88:ARG:HD2	4	0.11
(1,1687)	1:A:6:LEU:HD22	1:A:88:ARG:HD3	4	0.11
(1,1687)	1:A:6:LEU:HD23	1:A:88:ARG:HD2	4	0.11
(1,1687)	1:A:6:LEU:HD23	1:A:88:ARG:HD3	4	0.11
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	2	0.11
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	5	0.11
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	6	0.11
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	7	0.11
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	13	0.11
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	17	0.11
(1,1581)	1:A:34:ARG:HA	1:A:37:GLN:HE21	18	0.11
(1,1499)	1:A:49:GLN:HB3	1:A:98:LEU:H	19	0.11
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE2	2	0.11
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE3	2	0.11
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE2	2	0.11
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE3	2	0.11
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE2	3	0.11
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE3	3	0.11
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE2	3	0.11
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE3	3	0.11
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE2	20	0.11
(1,1207)	1:A:24:ASP:HB2	1:A:27:LYS:HE3	20	0.11
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE2	20	0.11
(1,1207)	1:A:24:ASP:HB3	1:A:27:LYS:HE3	20	0.11

## 10 Dihedral-angle violation analysis

Dihedral angle analysis failed due to data error in the dihedral angle restraints, possibly missing target value