



Full wwPDB NMR Structure Validation Report ⓘ

Mar 12, 2026 – 01:13 PM UTC

PDB ID : 6SJX / pdb_00006sjx
BMRB ID : 34424
Title : Precursor structure of the self-processing module of iron-regulated FrpC of N. Meningitidis with calcium ions
Authors : Kuban, V.; Macek, P.; Hritz, J.; Nechvatalova, K.; Nedbalcova, K.; Faldyna, M.; Zidek, L.; Bumba, L.
Deposited on : 2019-08-14

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

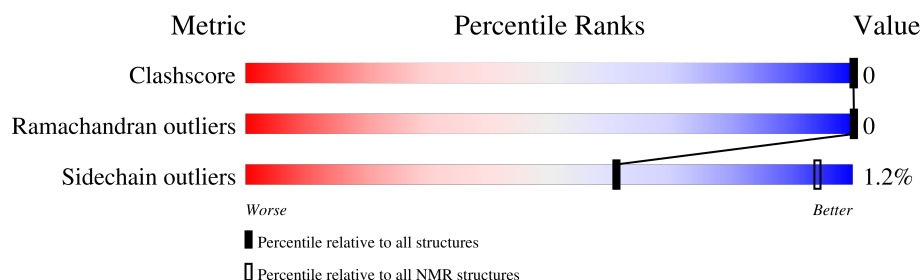
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 88%.

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	182	73% 27%

2 Ensemble composition and analysis

This entry contains 14 models. Model 1 is the overall representative, medoid model (most similar to other models).

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

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:0-A:15, A:36-A:49, A:55-A:66, A:78-A:168 (133)	0.99	1

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

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

The models can be grouped into 2 clusters and 2 single-model clusters were found.

Cluster number	Models
1	3, 4, 10, 11, 12, 13
2	1, 2, 5, 6, 8, 9
Single-model clusters	7; 14

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2651 atoms, of which 1291 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Iron-regulated protein FrpC.

Mol	Chain	Residues	Atoms						Trace
1	A	182	Total	C	H	N	O	S	0
			2646	829	1291	238	287	1	

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P55127
A	-1	SER	-	expression tag	UNP P55127
A	0	ASP	-	expression tag	UNP P55127
A	1	ALA	-	expression tag	UNP P55127
A	39	SER	ALA	conflict	UNP P55127
A	101	SER	THR	conflict	UNP P55127
A	157	ASN	ASP	conflict	UNP P55127
A	178	LEU	-	expression tag	UNP P55127
A	179	GLU	-	expression tag	UNP P55127

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

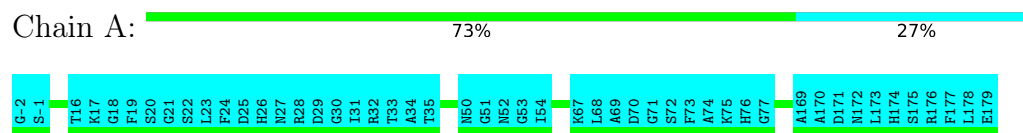
Mol	Chain	Residues	Atoms	
2	A	5	Total	Ca
			5	5

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: Iron-regulated protein FrpC

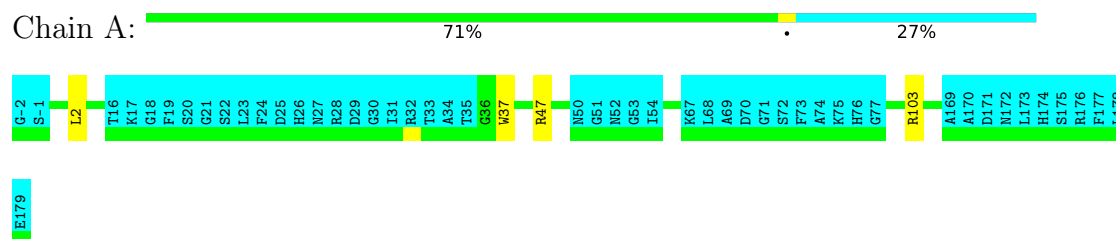


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

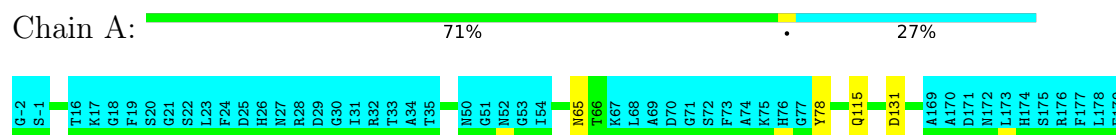
4.2.1 Score per residue for model 1 (medoid)

- Molecule 1: Iron-regulated protein FrpC



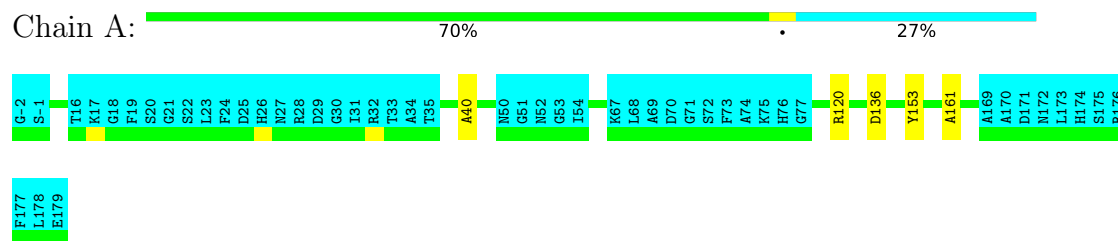
4.2.2 Score per residue for model 2

- Molecule 1: Iron-regulated protein FrpC



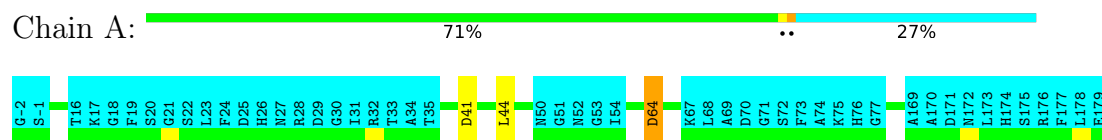
4.2.3 Score per residue for model 3

- Molecule 1: Iron-regulated protein FrpC



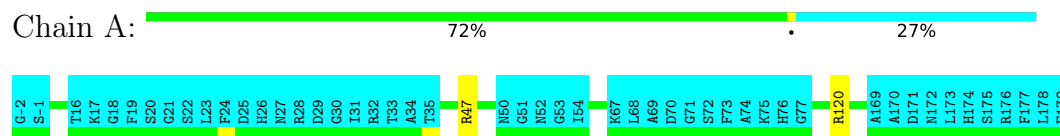
4.2.4 Score per residue for model 4

- Molecule 1: Iron-regulated protein FrpC



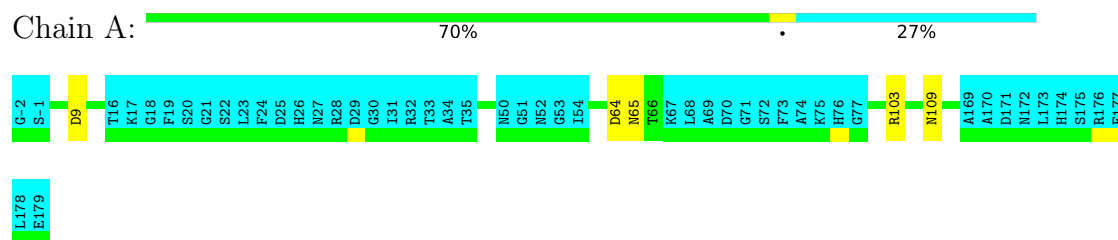
4.2.5 Score per residue for model 5

- Molecule 1: Iron-regulated protein FrpC



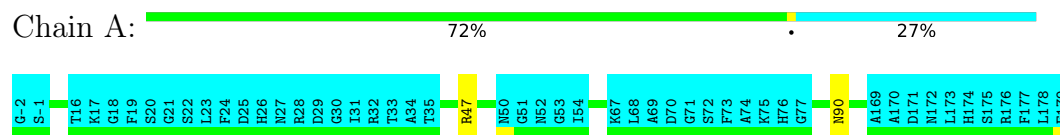
4.2.6 Score per residue for model 6

- Molecule 1: Iron-regulated protein FrpC



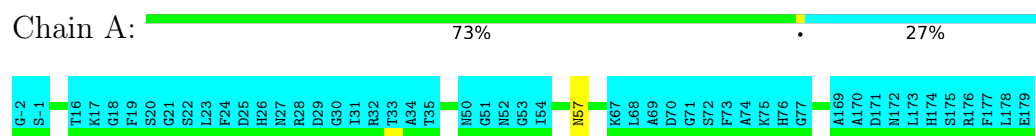
4.2.7 Score per residue for model 7

- Molecule 1: Iron-regulated protein FrpC



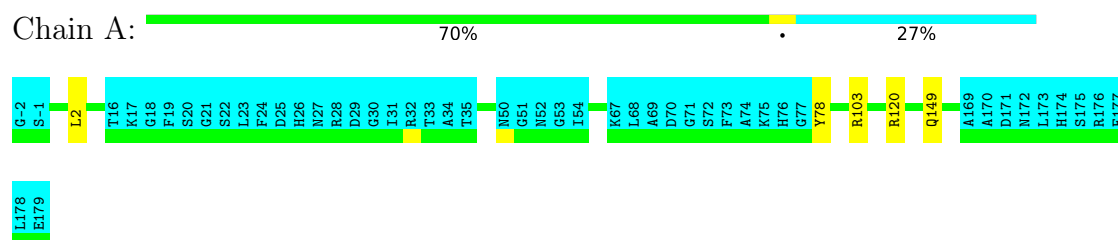
4.2.8 Score per residue for model 8

- Molecule 1: Iron-regulated protein FrpC



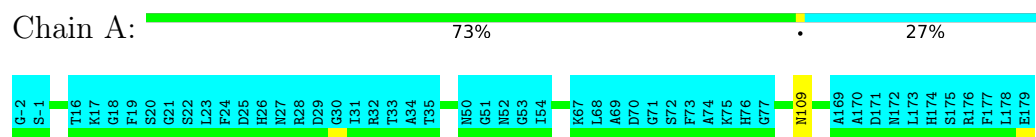
4.2.9 Score per residue for model 9

- Molecule 1: Iron-regulated protein FrpC



4.2.10 Score per residue for model 10

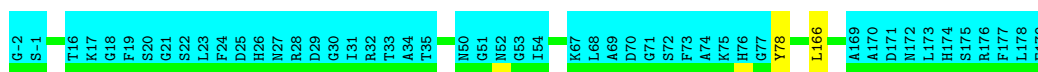
- Molecule 1: Iron-regulated protein FrpC



4.2.11 Score per residue for model 11

- Molecule 1: Iron-regulated protein FrpC

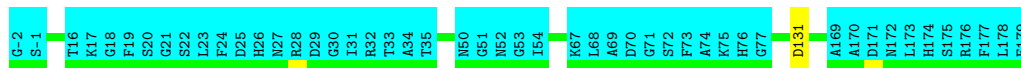




4.2.12 Score per residue for model 12

- Molecule 1: Iron-regulated protein FrpC

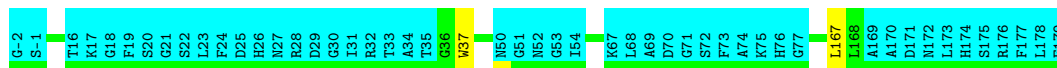
Chain A: 73% 27%



4.2.13 Score per residue for model 13

- Molecule 1: Iron-regulated protein FrpC

Chain A: 72% 27%



4.2.14 Score per residue for model 14

- Molecule 1: Iron-regulated protein FrpC

Chain A: 73% 27%



5 Refinement protocol and experimental data overview

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

Of the 20 calculated structures, 14 were deposited, based on the following criterion: *structures with the least restraint violations*.

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

Software name	Classification	Version
CYANA	structure calculation	3.97
GROMACS	refinement	5.1.1

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	1665
Number of shifts mapped to atoms	1665
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	88%

6 Model quality [i](#)

6.1 Standard geometry [i](#)

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

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	0.52±0.01	0±0/1000 (0.0± 0.0%)	1.39±0.02	2±1/1360 (0.1± 0.1%)
All	All	0.52	0/14000 (0.0%)	1.39	22/19040 (0.1%)

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	64	ASP	CA-CB-CG	6.94	119.54	112.60	4	2
1	A	47	ARG	NE-CZ-NH2	6.44	124.99	119.20	7	3
1	A	120	ARG	NE-CZ-NH2	6.43	124.99	119.20	9	3
1	A	131	ASP	CA-CB-CG	6.17	118.78	112.60	2	2
1	A	9	ASP	CA-CB-CG	5.70	118.30	112.60	6	2
1	A	57	ASN	CA-CB-CG	5.57	118.17	112.60	8	1
1	A	103	ARG	NE-CZ-NH2	5.48	124.13	119.20	9	3
1	A	136	ASP	CA-CB-CG	5.37	117.97	112.60	3	1
1	A	40	ALA	CA-C-N	5.27	127.86	120.28	3	1
1	A	40	ALA	C-N-CA	5.27	127.86	120.28	3	1
1	A	57	ASN	CB-CA-C	5.22	118.14	109.53	8	1
1	A	41	ASP	CA-C-N	5.11	130.24	122.11	4	1
1	A	41	ASP	C-N-CA	5.11	130.24	122.11	4	1

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	991	947	947	0±0
All	All	13944	13258	13258	1

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 0.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:153:TYR:CZ	1:A:161:ALA:HB3	0.40	2.52	3	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	133/182 (73%)	130±1 (98±1%)	3±1 (2±1%)	0±0 (0±0%)	100	100
All	All	1862/2548 (73%)	1820 (98%)	42 (2%)	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	103/139 (74%)	102±1 (99±1%)	1±1 (1±1%)	61	94
All	All	1442/1946 (74%)	1424 (99%)	18 (1%)	61	94

All 12 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	78	TYR	3
1	A	2	LEU	2
1	A	37	TRP	2
1	A	65	ASN	2
1	A	109	ASN	2
1	A	115	GLN	1
1	A	44	LEU	1
1	A	64	ASP	1
1	A	90	ASN	1
1	A	149	GLN	1
1	A	166	LEU	1
1	A	167	LEU	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

6.7 Other polymers ⓘ

There are no such molecules in this entry.

6.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 88% for the well-defined parts and 73% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *NMRStar-CS-spm-p1a.cif*

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	1665
Number of shifts mapped to atoms	1665
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	6

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	147	-0.14 ± 0.15	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	133	-0.06 ± 0.13	None needed (< 0.5 ppm)
$^{13}\text{C}'$	141	-0.01 ± 0.13	None needed (< 0.5 ppm)
^{15}N	144	0.54 ± 0.38	None needed (imprecise)

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 88%, i.e. 1479 atoms were assigned a chemical shift out of a possible 1682. 0 out of 26 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	644/679 (95%)	264/280 (94%)	254/266 (95%)	126/133 (95%)
Sidechain	812/932 (87%)	554/604 (92%)	247/296 (83%)	11/32 (34%)

Continued on next page...

Continued from previous page...

	Total	¹ H	¹³ C	¹⁵ N
Aromatic	23/71 (32%)	12/34 (35%)	10/35 (29%)	1/2 (50%)
Overall	1479/1682 (88%)	830/918 (90%)	511/597 (86%)	138/167 (83%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 73%, i.e. 1665 atoms were assigned a chemical shift out of a possible 2292. 0 out of 30 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	735/932 (79%)	303/386 (78%)	288/364 (79%)	144/182 (79%)
Sidechain	899/1228 (73%)	614/794 (77%)	274/386 (71%)	11/48 (23%)
Aromatic	31/132 (23%)	16/66 (24%)	14/61 (23%)	1/5 (20%)
Overall	1665/2292 (73%)	933/1246 (75%)	576/811 (71%)	156/235 (66%)

7.1.4 Statistically unusual chemical shifts ⓘ

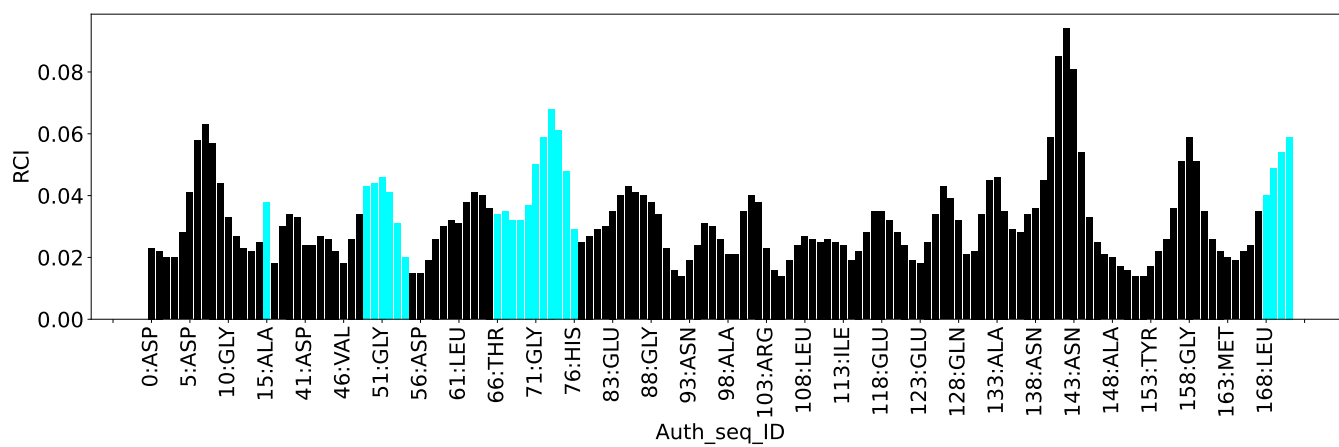
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	114	SER	HB2	1.49	2.61 – 5.13	-9.4
1	A	118	GLU	HB3	3.33	0.95 – 3.05	6.3
1	A	44	LEU	HD11	-0.62	-0.61 – 2.12	-5.0
1	A	44	LEU	HD12	-0.62	-0.61 – 2.12	-5.0
1	A	44	LEU	HD13	-0.62	-0.61 – 2.12	-5.0
1	A	107	ASP	HB2	1.40	1.41 – 4.01	-5.0

7.1.5 Random Coil Index (RCI) plots ⓘ

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	2056
Intra-residue ($ i-j =0$)	512
Sequential ($ i-j =1$)	506
Medium range ($ i-j >1$ and $ i-j <5$)	278
Long range ($ i-j \geq 5$)	760
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	238
Number of unmapped restraints	0
Number of restraints per residue	12.3
Number of long range restraints per residue ¹	4.1

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)	4.8	0.2
0.2-0.5 (Medium)	7.0	0.5
>0.5 (Large)	2.2	1.16

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)	9.1	6.8
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

9 Distance violation analysis [i](#)

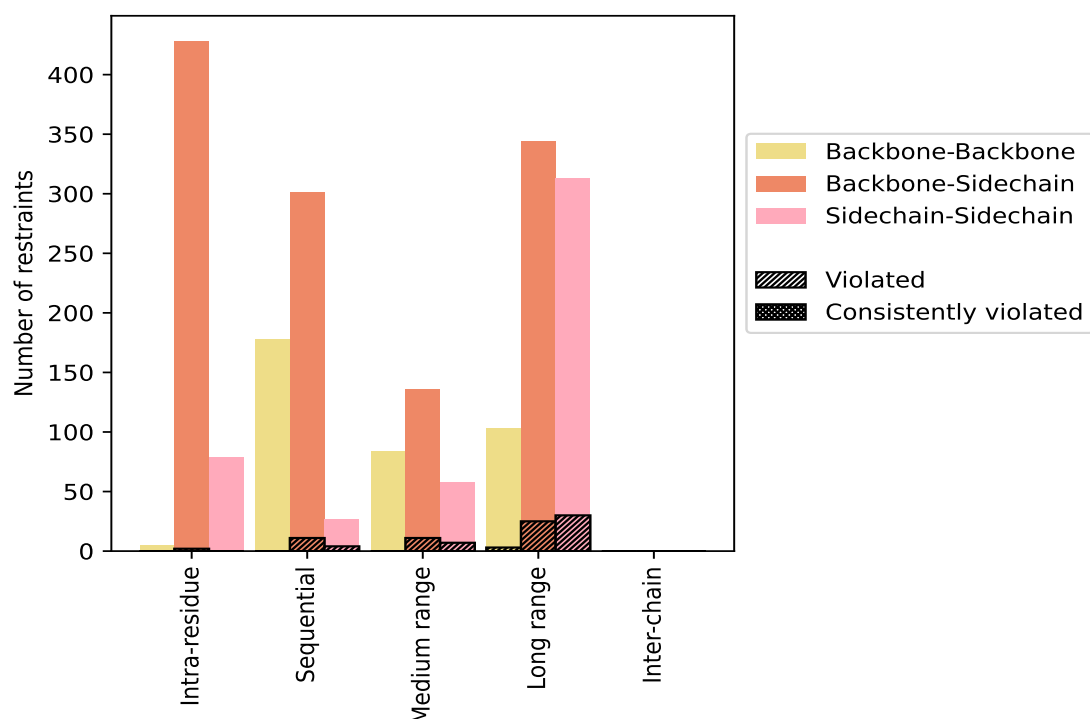
9.1 Summary of distance violations [i](#)

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

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	512	24.9	2	0.4	0.1	0	0.0	0.0
Backbone-Backbone	5	0.2	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	428	20.8	2	0.5	0.1	0	0.0	0.0
Sidechain-Sidechain	79	3.8	0	0.0	0.0	0	0.0	0.0
Sequential ($i-j =1$)	506	24.6	15	3.0	0.7	0	0.0	0.0
Backbone-Backbone	178	8.7	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	301	14.6	11	3.7	0.5	0	0.0	0.0
Sidechain-Sidechain	27	1.3	4	14.8	0.2	0	0.0	0.0
Medium range ($i-j >1$ & $i-j <5$)	278	13.5	18	6.5	0.9	0	0.0	0.0
Backbone-Backbone	84	4.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	136	6.6	11	8.1	0.5	0	0.0	0.0
Sidechain-Sidechain	58	2.8	7	12.1	0.3	0	0.0	0.0
Long range ($i-j \geq 5$)	760	37.0	58	7.6	2.8	0	0.0	0.0
Backbone-Backbone	103	5.0	3	2.9	0.1	0	0.0	0.0
Backbone-Sidechain	344	16.7	25	7.3	1.2	0	0.0	0.0
Sidechain-Sidechain	313	15.2	30	9.6	1.5	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	2056	100.0	93	4.5	4.5	0	0.0	0.0
Backbone-Backbone	370	18.0	3	0.8	0.1	0	0.0	0.0
Backbone-Sidechain	1209	58.8	49	4.1	2.4	0	0.0	0.0
Sidechain-Sidechain	477	23.2	41	8.6	2.0	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	5	2	9	0	16	0.29	0.8	0.18	0.2
2	0	3	1	2	0	6	0.28	0.73	0.22	0.17
3	2	3	7	7	0	19	0.33	1.0	0.21	0.27
4	0	3	2	8	0	13	0.38	0.93	0.23	0.27
5	0	2	9	4	0	15	0.25	0.67	0.15	0.21
6	1	2	0	10	0	13	0.32	0.59	0.14	0.27
7	0	3	3	14	0	20	0.33	0.79	0.19	0.26
8	1	2	3	10	0	16	0.35	0.86	0.24	0.29
9	0	3	2	5	0	10	0.32	0.93	0.23	0.26
10	0	1	2	5	0	8	0.31	0.92	0.27	0.15

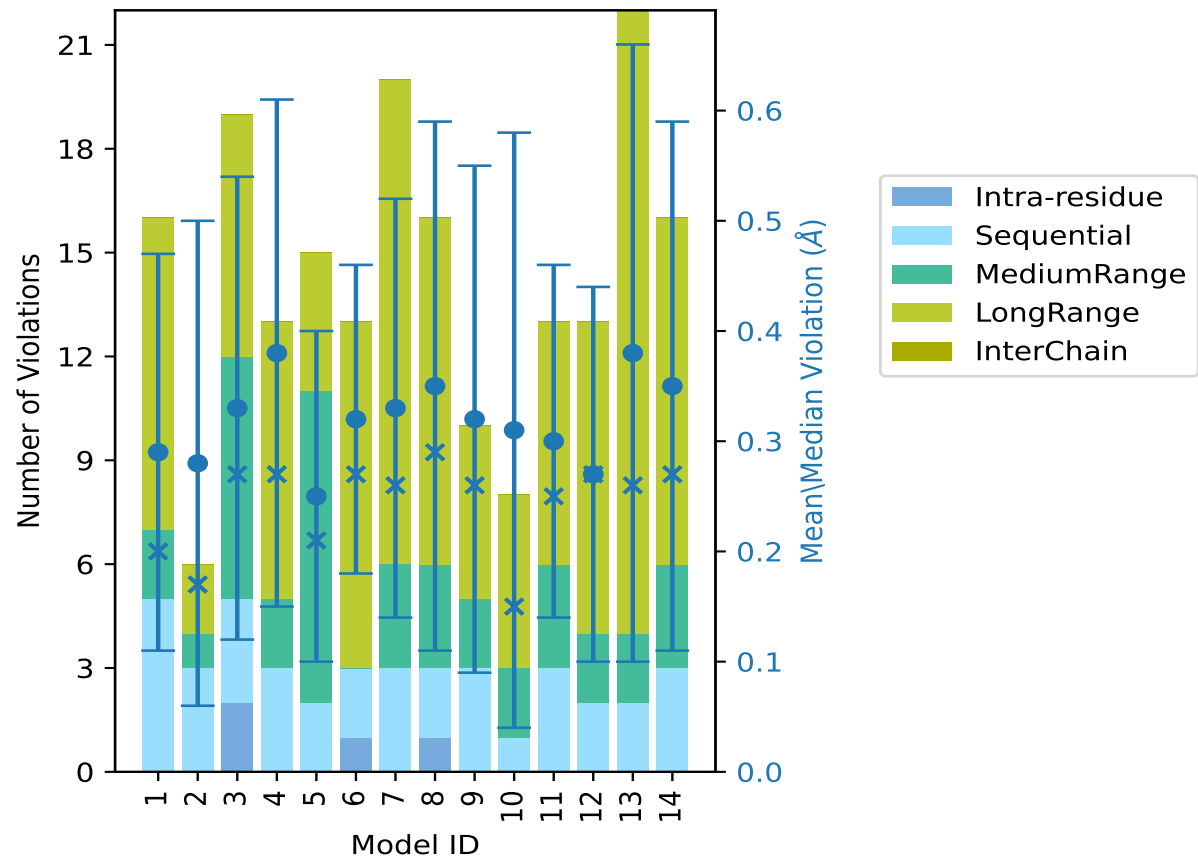
Continued on next page...

Continued from previous page...

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	0	3	3	7	0	13	0.3	0.66	0.16	0.25
12	0	2	2	9	0	13	0.27	0.75	0.17	0.27
13	0	2	2	18	0	22	0.38	1.16	0.28	0.26
14	0	3	3	10	0	16	0.35	1.09	0.24	0.27

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

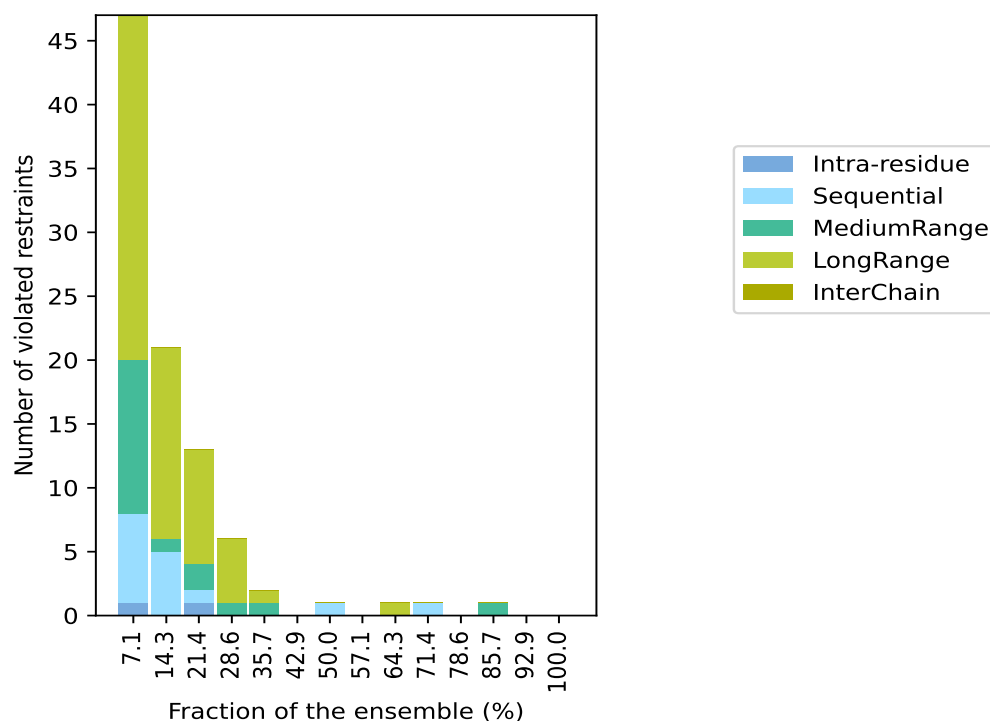
a given fraction of the ensemble. In total, 1963(IR:510, SQ:491, MR:260, LR:702, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
1	7	12	27	0	47	1	7.1
0	5	1	15	0	21	2	14.3
1	1	2	9	0	13	3	21.4
0	0	1	5	0	6	4	28.6
0	0	1	1	0	2	5	35.7
0	0	0	0	0	0	6	42.9
0	1	0	0	0	1	7	50.0
0	0	0	0	0	0	8	57.1
0	0	0	1	0	1	9	64.3
0	1	0	0	0	1	10	71.4
0	0	0	0	0	0	11	78.6
0	0	1	0	0	1	12	85.7
0	0	0	0	0	0	13	92.9
0	0	0	0	0	0	14	100.0

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

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

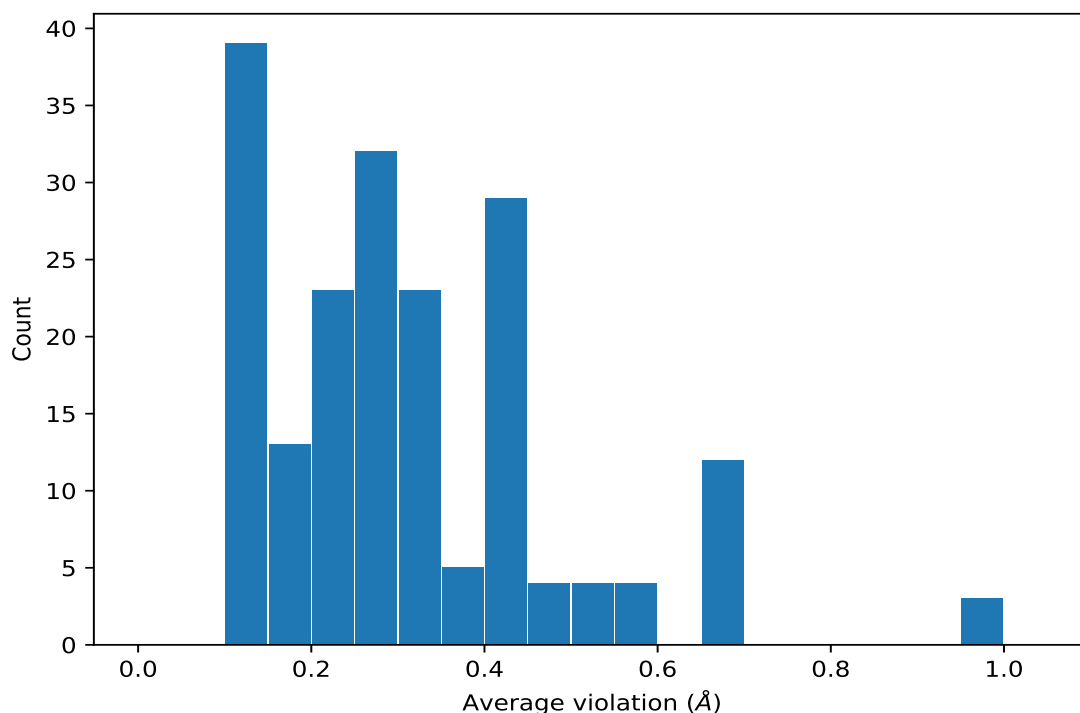
9.3.1 Bar graph : Distance violation statistics for the ensemble ⓘ



9.4 Most violated distance restraints in the ensemble ⓘ

9.4.1 Histogram : Distribution of mean distance violations ⓘ

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



9.4.2 Table: Most violated distance restraints ⓘ

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)
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	12	0.6	0.15	0.63
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	12	0.6	0.15	0.63
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	12	0.6	0.15	0.63
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	10	0.31	0.03	0.32
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	10	0.31	0.03	0.32
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	10	0.31	0.03	0.32
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	10	0.31	0.03	0.32
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	10	0.31	0.03	0.32
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	10	0.31	0.03	0.32

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	9	0.23	0.07	0.23
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	9	0.23	0.07	0.23
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD21	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD22	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD21	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD22	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD21	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD22	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD21	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD22	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD21	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD22	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD21	7	0.31	0.14	0.28
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD22	7	0.31	0.14	0.28
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD21	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD22	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD21	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD22	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD21	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD22	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD21	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD22	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD21	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD22	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD21	5	0.65	0.3	0.69
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD22	5	0.65	0.3	0.69
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD11	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD12	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD13	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD21	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD22	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD23	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD11	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD12	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD13	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD21	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD22	5	0.26	0.07	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD23	5	0.26	0.07	0.25
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG2	4	0.54	0.33	0.55
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG3	4	0.54	0.33	0.55
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE1	4	0.46	0.12	0.44
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE2	4	0.46	0.12	0.44

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE3	4	0.46	0.12	0.44
(1,93)	1:40:A:ALA:HB1	1:108:A:LEU:HA	4	0.43	0.25	0.34
(1,93)	1:40:A:ALA:HB2	1:108:A:LEU:HA	4	0.43	0.25	0.34
(1,93)	1:40:A:ALA:HB3	1:108:A:LEU:HA	4	0.43	0.25	0.34
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	4	0.38	0.04	0.4
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	4	0.38	0.04	0.4
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	4	0.38	0.04	0.4
(2,137)	1:9:A:ASP:H	1:162:A:LYS:H	4	0.32	0.12	0.34
(2,75)	1:135:A:LYS:HE2	1:137:A:VAL:HB	4	0.3	0.13	0.29
(2,75)	1:135:A:LYS:HE3	1:137:A:VAL:HB	4	0.3	0.13	0.29
(1,134)	1:147:A:LEU:HD21	1:166:A:LEU:HG	3	0.98	0.13	0.93
(1,134)	1:147:A:LEU:HD22	1:166:A:LEU:HG	3	0.98	0.13	0.93
(1,134)	1:147:A:LEU:HD23	1:166:A:LEU:HG	3	0.98	0.13	0.93
(1,484)	1:135:A:LYS:HG2	1:136:A:ASP:HA	3	0.45	0.07	0.47
(1,555)	1:134:A:TYR:HE1	1:149:A:GLN:HG2	3	0.42	0.11	0.47
(1,555)	1:134:A:TYR:HE2	1:149:A:GLN:HG2	3	0.42	0.11	0.47
(1,1643)	1:108:A:LEU:HD11	1:118:A:GLU:HA	3	0.4	0.13	0.48
(1,1643)	1:108:A:LEU:HD12	1:118:A:GLU:HA	3	0.4	0.13	0.48
(1,1643)	1:108:A:LEU:HD13	1:118:A:GLU:HA	3	0.4	0.13	0.48
(1,1643)	1:108:A:LEU:HD21	1:118:A:GLU:HA	3	0.4	0.13	0.48
(1,1643)	1:108:A:LEU:HD22	1:118:A:GLU:HA	3	0.4	0.13	0.48
(1,1643)	1:108:A:LEU:HD23	1:118:A:GLU:HA	3	0.4	0.13	0.48
(1,1753)	1:134:A:TYR:HE1	1:149:A:GLN:HG2	3	0.35	0.08	0.34
(1,1753)	1:134:A:TYR:HE1	1:149:A:GLN:HG3	3	0.35	0.08	0.34
(1,1753)	1:134:A:TYR:HE2	1:149:A:GLN:HG2	3	0.35	0.08	0.34
(1,1753)	1:134:A:TYR:HE2	1:149:A:GLN:HG3	3	0.35	0.08	0.34
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG21	3	0.27	0.14	0.24
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG22	3	0.27	0.14	0.24
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG23	3	0.27	0.14	0.24
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB1	3	0.26	0.11	0.29
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB2	3	0.26	0.11	0.29
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB3	3	0.26	0.11	0.29
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB1	3	0.26	0.11	0.29
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB2	3	0.26	0.11	0.29
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB3	3	0.26	0.11	0.29
(1,272)	1:94:A:ALA:HA	1:100:A:GLN:HG2	3	0.22	0.11	0.17
(1,272)	1:94:A:ALA:HA	1:100:A:GLN:HG3	3	0.22	0.11	0.17
(1,218)	1:68:A:LEU:HB3	1:83:A:GLU:HG2	3	0.18	0.05	0.21
(1,622)	1:46:A:VAL:HA	1:61:A:LEU:HB2	3	0.18	0.05	0.17
(1,1420)	1:47:A:ARG:HB2	1:49:A:LEU:H	3	0.17	0.02	0.18
(1,1420)	1:47:A:ARG:HB3	1:49:A:LEU:H	3	0.17	0.02	0.18
(1,736)	1:61:A:LEU:H	1:61:A:LEU:HG	3	0.15	0.03	0.15

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB1	3	0.14	0.0	0.14
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB2	3	0.14	0.0	0.14
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB3	3	0.14	0.0	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	3	0.14	0.0	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	3	0.14	0.0	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	3	0.14	0.0	0.14
(1,255)	1:107:A:ASP:HA	1:108:A:LEU:HG	2	0.55	0.11	0.55
(1,1724)	1:128:A:GLN:HE21	1:156:A:THR:HB	2	0.51	0.08	0.51
(1,1724)	1:128:A:GLN:HE22	1:156:A:THR:HB	2	0.51	0.08	0.51
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD11	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD12	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD13	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD21	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD22	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD23	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD11	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD12	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD13	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD21	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD22	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD23	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD11	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD12	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD13	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD21	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD22	2	0.4	0.02	0.4
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD23	2	0.4	0.02	0.4
(1,1650)	1:114:A:SER:HB2	1:119:A:LEU:HG	2	0.38	0.15	0.38
(1,1650)	1:114:A:SER:HB3	1:119:A:LEU:HG	2	0.38	0.15	0.38
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD21	2	0.29	0.11	0.29
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD22	2	0.29	0.11	0.29
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD23	2	0.29	0.11	0.29
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD21	2	0.29	0.11	0.29
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD22	2	0.29	0.11	0.29
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD23	2	0.29	0.11	0.29
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD21	2	0.29	0.11	0.29
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD22	2	0.29	0.11	0.29
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD23	2	0.29	0.11	0.29
(1,519)	1:163:A:MET:HE1	1:164:A:GLY:HA2	2	0.24	0.05	0.24
(1,519)	1:163:A:MET:HE2	1:164:A:GLY:HA2	2	0.24	0.05	0.24
(1,519)	1:163:A:MET:HE3	1:164:A:GLY:HA2	2	0.24	0.05	0.24
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD21	2	0.24	0.0	0.24

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD22	2	0.24	0.0	0.24
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD23	2	0.24	0.0	0.24
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD21	2	0.24	0.03	0.24
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD22	2	0.24	0.03	0.24
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD23	2	0.24	0.03	0.24
(1,62)	1:11:A:ILE:HD11	1:162:A:LYS:HA	2	0.22	0.09	0.22
(1,62)	1:11:A:ILE:HD12	1:162:A:LYS:HA	2	0.22	0.09	0.22
(1,62)	1:11:A:ILE:HD13	1:162:A:LYS:HA	2	0.22	0.09	0.22
(2,136)	1:2:A:LEU:HG	1:166:A:LEU:H	2	0.2	0.03	0.2
(2,186)	1:81:A:LEU:HD11	1:130:A:LEU:H	2	0.2	0.05	0.2
(2,186)	1:81:A:LEU:HD12	1:130:A:LEU:H	2	0.2	0.05	0.2
(2,186)	1:81:A:LEU:HD13	1:130:A:LEU:H	2	0.2	0.05	0.2
(2,186)	1:81:A:LEU:HD21	1:130:A:LEU:H	2	0.2	0.05	0.2
(2,186)	1:81:A:LEU:HD22	1:130:A:LEU:H	2	0.2	0.05	0.2
(2,186)	1:81:A:LEU:HD23	1:130:A:LEU:H	2	0.2	0.05	0.2
(1,210)	1:59:A:ALA:HA	1:66:A:THR:HA	2	0.18	0.08	0.18
(2,198)	1:103:A:ARG:HD2	1:122:A:LEU:H	2	0.18	0.04	0.18
(2,198)	1:103:A:ARG:HD3	1:122:A:LEU:H	2	0.18	0.04	0.18
(2,76)	1:135:A:LYS:HE2	1:150:A:GLN:HB2	2	0.18	0.07	0.18
(2,76)	1:135:A:LYS:HE2	1:150:A:GLN:HB3	2	0.18	0.07	0.18
(2,76)	1:135:A:LYS:HE3	1:150:A:GLN:HB2	2	0.18	0.07	0.18
(2,76)	1:135:A:LYS:HE3	1:150:A:GLN:HB3	2	0.18	0.07	0.18
(1,1566)	1:90:A:ASN:HB3	1:91:A:ILE:HG12	2	0.16	0.03	0.16
(1,1566)	1:90:A:ASN:HB3	1:91:A:ILE:HG13	2	0.16	0.03	0.16
(1,1360)	1:11:A:ILE:HG12	1:163:A:MET:HB3	2	0.15	0.01	0.15
(1,1360)	1:11:A:ILE:HG13	1:163:A:MET:HB3	2	0.15	0.01	0.15
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD11	2	0.15	0.04	0.15
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD12	2	0.15	0.04	0.15
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD13	2	0.15	0.04	0.15
(2,68)	1:0:A:ASP:HA	1:166:A:LEU:HB2	2	0.13	0.01	0.13
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE1	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE2	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE3	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE1	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE2	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE3	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE1	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE2	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE3	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE1	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE2	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE3	2	0.12	0.02	0.12

Continued on next page...

Continued from previous page...

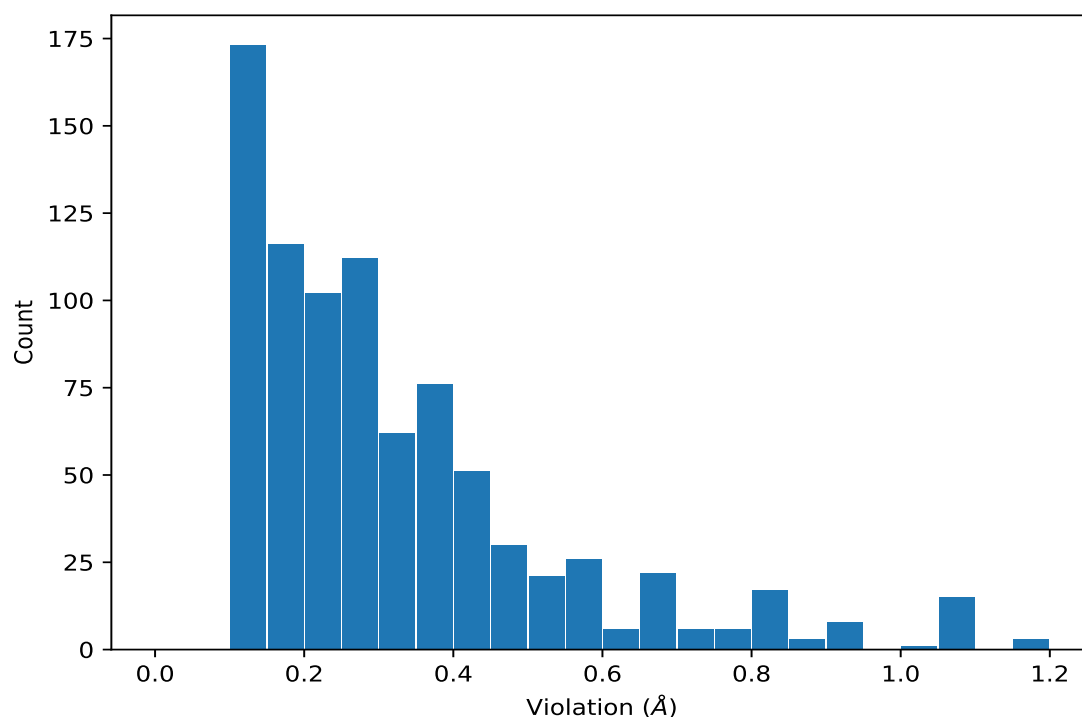
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE1	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE2	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE3	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE1	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE2	2	0.12	0.02	0.12
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE3	2	0.12	0.02	0.12
(2,187)	1:82:A:ALA:H	1:83:A:GLU:HG2	2	0.12	0.0	0.12
(2,187)	1:82:A:ALA:H	1:83:A:GLU:HG3	2	0.12	0.0	0.12
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD11	2	0.12	0.02	0.12
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD12	2	0.12	0.02	0.12
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD13	2	0.12	0.02	0.12
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD21	2	0.12	0.02	0.12
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD22	2	0.12	0.02	0.12
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD23	2	0.12	0.02	0.12

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

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)
(1,134)	1:147:A:LEU:HD21	1:166:A:LEU:HG	13	1.16
(1,134)	1:147:A:LEU:HD22	1:166:A:LEU:HG	13	1.16
(1,134)	1:147:A:LEU:HD23	1:166:A:LEU:HG	13	1.16
(1,145)	1:48:A:ASP:HA	1:55:A:ILE:HD11	14	1.09
(1,145)	1:48:A:ASP:HA	1:55:A:ILE:HD12	14	1.09
(1,145)	1:48:A:ASP:HA	1:55:A:ILE:HD13	14	1.09
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD21	13	1.05
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD22	13	1.05
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD21	13	1.05
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD22	13	1.05
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD21	13	1.05
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD22	13	1.05
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD21	13	1.05
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD22	13	1.05
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD21	13	1.05
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD22	13	1.05
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD21	13	1.05
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD22	13	1.05
(1,915)	1:101:A:SER:H	1:102:A:LEU:HG	3	1.0
(1,343)	1:90:A:ASN:HB3	1:91:A:ILE:HG21	4	0.93
(1,343)	1:90:A:ASN:HB3	1:91:A:ILE:HG22	4	0.93
(1,343)	1:90:A:ASN:HB3	1:91:A:ILE:HG23	4	0.93
(1,134)	1:147:A:LEU:HD21	1:166:A:LEU:HG	9	0.93
(1,134)	1:147:A:LEU:HD22	1:166:A:LEU:HG	9	0.93
(1,134)	1:147:A:LEU:HD23	1:166:A:LEU:HG	9	0.93
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG2	10	0.92
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG3	10	0.92
(1,134)	1:147:A:LEU:HD21	1:166:A:LEU:HG	8	0.86
(1,134)	1:147:A:LEU:HD22	1:166:A:LEU:HG	8	0.86
(1,134)	1:147:A:LEU:HD23	1:166:A:LEU:HG	8	0.86
(1,93)	1:40:A:ALA:HB1	1:108:A:LEU:HA	13	0.83
(1,93)	1:40:A:ALA:HB2	1:108:A:LEU:HA	13	0.83
(1,93)	1:40:A:ALA:HB3	1:108:A:LEU:HA	13	0.83
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG2	8	0.81
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG3	8	0.81
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD21	1	0.8
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD22	1	0.8

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD21	1	0.8
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD22	1	0.8
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD21	1	0.8
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD22	1	0.8
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD21	1	0.8
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD22	1	0.8
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD21	1	0.8
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD22	1	0.8
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD21	1	0.8
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD22	1	0.8
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	7	0.79
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	7	0.79
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	7	0.79
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	12	0.75
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	12	0.75
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	12	0.75
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	2	0.73
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	2	0.73
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	2	0.73
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	8	0.72
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	8	0.72
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	8	0.72
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD21	7	0.69
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD22	7	0.69
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD21	7	0.69
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD22	7	0.69
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD21	7	0.69
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD22	7	0.69
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD21	7	0.69
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD22	7	0.69
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD21	7	0.69
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD22	7	0.69
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD21	7	0.69
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD22	7	0.69
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	5	0.67
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	5	0.67
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	5	0.67
(1,255)	1:107:A:ASP:HA	1:108:A:LEU:HG	14	0.66
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	11	0.66
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	11	0.66
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	11	0.66
(1,1272)	1:91:A:ILE:HD11	1:93:A:ASN:HD22	4	0.65

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1272)	1:91:A:ILE:HD12	1:93:A:ASN:HD22	4	0.65
(1,1272)	1:91:A:ILE:HD13	1:93:A:ASN:HD22	4	0.65
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE1	7	0.64
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE2	7	0.64
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE3	7	0.64
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	3	0.6
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	3	0.6
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	3	0.6
(1,1724)	1:128:A:GLN:HE21	1:156:A:THR:HB	6	0.59
(1,1724)	1:128:A:GLN:HE22	1:156:A:THR:HB	6	0.59
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD21	4	0.58
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD22	4	0.58
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD21	4	0.58
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD22	4	0.58
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD21	4	0.58
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD22	4	0.58
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD21	4	0.58
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD22	4	0.58
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD21	4	0.58
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD22	4	0.58
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD21	4	0.58
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD22	4	0.58
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD21	1	0.55
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD22	1	0.55
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD21	1	0.55
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD22	1	0.55
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD21	1	0.55
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD22	1	0.55
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD21	1	0.55
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD22	1	0.55
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD21	1	0.55
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD22	1	0.55
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD21	1	0.55
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD22	1	0.55
(1,1145)	1:153:A:TYR:H	1:160:A:THR:HB	14	0.54
(1,1650)	1:114:A:SER:HB2	1:119:A:LEU:HG	13	0.52
(1,1650)	1:114:A:SER:HB3	1:119:A:LEU:HG	13	0.52
(1,555)	1:134:A:TYR:HE1	1:149:A:GLN:HG2	3	0.52
(1,555)	1:134:A:TYR:HE2	1:149:A:GLN:HG2	3	0.52
(1,484)	1:135:A:LYS:HG2	1:136:A:ASP:HA	4	0.52
(1,1643)	1:108:A:LEU:HD11	1:118:A:GLU:HA	7	0.51
(1,1643)	1:108:A:LEU:HD12	1:118:A:GLU:HA	7	0.51

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1643)	1:108:A:LEU:HD13	1:118:A:GLU:HA	7	0.51
(1,1643)	1:108:A:LEU:HD21	1:118:A:GLU:HA	7	0.51
(1,1643)	1:108:A:LEU:HD22	1:118:A:GLU:HA	7	0.51
(1,1643)	1:108:A:LEU:HD23	1:118:A:GLU:HA	7	0.51
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	9	0.51
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	9	0.51
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	9	0.51
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE1	13	0.5
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE2	13	0.5
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE3	13	0.5
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	13	0.5
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	13	0.5
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	13	0.5
(1,1643)	1:108:A:LEU:HD11	1:118:A:GLU:HA	6	0.48
(1,1643)	1:108:A:LEU:HD12	1:118:A:GLU:HA	6	0.48
(1,1643)	1:108:A:LEU:HD13	1:118:A:GLU:HA	6	0.48
(1,1643)	1:108:A:LEU:HD21	1:118:A:GLU:HA	6	0.48
(1,1643)	1:108:A:LEU:HD22	1:118:A:GLU:HA	6	0.48
(1,1643)	1:108:A:LEU:HD23	1:118:A:GLU:HA	6	0.48
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	10	0.48
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	10	0.48
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	10	0.48
(2,137)	1:9:A:ASP:H	1:162:A:LYS:H	6	0.47
(1,1463)	1:67:A:LYS:HD2	1:71:A:GLY:H	14	0.47
(1,1463)	1:67:A:LYS:HD3	1:71:A:GLY:H	14	0.47
(1,555)	1:134:A:TYR:HE1	1:149:A:GLN:HG2	8	0.47
(1,555)	1:134:A:TYR:HE2	1:149:A:GLN:HG2	8	0.47
(1,484)	1:135:A:LYS:HG2	1:136:A:ASP:HA	5	0.47
(2,75)	1:135:A:LYS:HE2	1:137:A:VAL:HB	3	0.46
(2,75)	1:135:A:LYS:HE3	1:137:A:VAL:HB	3	0.46
(1,1753)	1:134:A:TYR:HE1	1:149:A:GLN:HG2	3	0.46
(1,1753)	1:134:A:TYR:HE1	1:149:A:GLN:HG3	3	0.46
(1,1753)	1:134:A:TYR:HE2	1:149:A:GLN:HG2	3	0.46
(1,1753)	1:134:A:TYR:HE2	1:149:A:GLN:HG3	3	0.46
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	1	0.46
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	1	0.46
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	1	0.46
(1,93)	1:40:A:ALA:HB1	1:108:A:LEU:HA	11	0.46
(1,93)	1:40:A:ALA:HB2	1:108:A:LEU:HA	11	0.46
(1,93)	1:40:A:ALA:HB3	1:108:A:LEU:HA	11	0.46
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG21	3	0.46
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG22	3	0.46

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG23	3	0.46
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD21	6	0.44
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD22	6	0.44
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD21	6	0.44
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD22	6	0.44
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD21	6	0.44
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD22	6	0.44
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD21	6	0.44
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD22	6	0.44
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD21	6	0.44
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD22	6	0.44
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD21	6	0.44
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD22	6	0.44
(1,255)	1:107:A:ASP:HA	1:108:A:LEU:HG	10	0.44
(1,1724)	1:128:A:GLN:HE21	1:156:A:THR:HB	11	0.43
(1,1724)	1:128:A:GLN:HE22	1:156:A:THR:HB	11	0.43
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD11	11	0.43
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD12	11	0.43
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD13	11	0.43
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD21	11	0.43
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD22	11	0.43
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD23	11	0.43
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD11	11	0.43
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD12	11	0.43
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD13	11	0.43
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD21	11	0.43
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD22	11	0.43
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD23	11	0.43
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD11	11	0.43
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD12	11	0.43
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD13	11	0.43
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD21	11	0.43
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD22	11	0.43
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD23	11	0.43
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	7	0.41
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	7	0.41
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	7	0.41
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD21	4	0.4
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD22	4	0.4
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD23	4	0.4
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD21	4	0.4
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD22	4	0.4

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD23	4	0.4
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD21	4	0.4
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD22	4	0.4
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD23	4	0.4
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	1	0.4
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	1	0.4
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	1	0.4
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	11	0.4
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	11	0.4
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	11	0.4
(2,75)	1:135:A:LYS:HE2	1:137:A:VAL:HB	7	0.39
(2,75)	1:135:A:LYS:HE3	1:137:A:VAL:HB	7	0.39
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE1	8	0.39
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE2	8	0.39
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE3	8	0.39
(1,1808)	1:152:A:SER:HB2	1:160:A:THR:HG21	4	0.39
(1,1808)	1:152:A:SER:HB2	1:160:A:THR:HG22	4	0.39
(1,1808)	1:152:A:SER:HB2	1:160:A:THR:HG23	4	0.39
(1,1808)	1:152:A:SER:HB3	1:160:A:THR:HG21	4	0.39
(1,1808)	1:152:A:SER:HB3	1:160:A:THR:HG22	4	0.39
(1,1808)	1:152:A:SER:HB3	1:160:A:THR:HG23	4	0.39
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD21	8	0.38
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD22	8	0.38
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD21	8	0.38
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD22	8	0.38
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD21	8	0.38
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD22	8	0.38
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD21	8	0.38
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD22	8	0.38
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD21	8	0.38
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD22	8	0.38
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD21	8	0.38
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD22	8	0.38
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	1	0.38
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	1	0.38
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB1	9	0.38
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB2	9	0.38
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB3	9	0.38
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB1	9	0.38
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB2	9	0.38
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB3	9	0.38
(1,1816)	1:161:A:ALA:HA	1:162:A:LYS:HD2	14	0.38

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1816)	1:161:A:ALA:HA	1:162:A:LYS:HD3	14	0.38
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD11	13	0.38
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD12	13	0.38
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD13	13	0.38
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD21	13	0.38
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD22	13	0.38
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD23	13	0.38
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD11	13	0.38
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD12	13	0.38
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD13	13	0.38
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD21	13	0.38
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD22	13	0.38
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD23	13	0.38
(1,1461)	1:67:A:LYS:HG2	1:71:A:GLY:H	14	0.38
(1,1461)	1:67:A:LYS:HG3	1:71:A:GLY:H	14	0.38
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD11	13	0.38
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD12	13	0.38
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD13	13	0.38
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD21	13	0.38
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD22	13	0.38
(1,1392)	1:40:A:ALA:HB1	1:108:A:LEU:HD23	13	0.38
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD11	13	0.38
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD12	13	0.38
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD13	13	0.38
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD21	13	0.38
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD22	13	0.38
(1,1392)	1:40:A:ALA:HB2	1:108:A:LEU:HD23	13	0.38
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD11	13	0.38
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD12	13	0.38
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD13	13	0.38
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD21	13	0.38
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD22	13	0.38
(1,1392)	1:40:A:ALA:HB3	1:108:A:LEU:HD23	13	0.38
(1,292)	1:87:A:ASN:HD21	1:95:A:ALA:HB1	12	0.38
(1,292)	1:87:A:ASN:HD21	1:95:A:ALA:HB2	12	0.38
(1,292)	1:87:A:ASN:HD21	1:95:A:ALA:HB3	12	0.38
(1,272)	1:94:A:ALA:HA	1:100:A:GLN:HG2	7	0.37
(1,272)	1:94:A:ALA:HA	1:100:A:GLN:HG3	7	0.37
(1,1277)	1:0:A:ASP:HB2	1:167:A:LEU:HG	6	0.36
(1,1277)	1:0:A:ASP:HB3	1:167:A:LEU:HG	6	0.36
(1,484)	1:135:A:LYS:HG2	1:136:A:ASP:HA	7	0.35
(1,407)	1:91:A:ILE:HG21	1:129:A:SER:HB3	13	0.35

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,407)	1:91:A:ILE:HG22	1:129:A:SER:HB3	13	0.35
(1,407)	1:91:A:ILE:HG23	1:129:A:SER:HB3	13	0.35
(2,137)	1:9:A:ASP:H	1:162:A:LYS:H	12	0.34
(1,1753)	1:134:A:TYR:HE1	1:149:A:GLN:HG2	8	0.34
(1,1753)	1:134:A:TYR:HE1	1:149:A:GLN:HG3	8	0.34
(1,1753)	1:134:A:TYR:HE2	1:149:A:GLN:HG2	8	0.34
(1,1753)	1:134:A:TYR:HE2	1:149:A:GLN:HG3	8	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	2	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	2	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	2	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	2	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	2	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	2	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	3	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	3	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	3	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	3	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	3	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	3	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	8	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	8	0.34
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	8	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	8	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	8	0.34
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	8	0.34
(2,137)	1:9:A:ASP:H	1:162:A:LYS:H	14	0.33
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	1	0.33
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	1	0.33
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	1	0.33
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	1	0.33
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	1	0.33
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	1	0.33
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	11	0.33
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	11	0.33
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	11	0.33
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	11	0.33
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	11	0.33
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	11	0.33
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE1	12	0.31
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE2	12	0.31
(2,46)	1:3:A:ALA:H	1:163:A:MET:HE3	12	0.31
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	7	0.31

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	7	0.31
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	7	0.31
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	7	0.31
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	7	0.31
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	7	0.31
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	3	0.31
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	3	0.31
(1,290)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	3	0.31
(1,62)	1:11:A:ILE:HD11	1:162:A:LYS:HA	5	0.31
(1,62)	1:11:A:ILE:HD12	1:162:A:LYS:HA	5	0.31
(1,62)	1:11:A:ILE:HD13	1:162:A:LYS:HA	5	0.31
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	12	0.3
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	12	0.3
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	12	0.3
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	12	0.3
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	12	0.3
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	12	0.3
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	12	0.3
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	12	0.3
(1,519)	1:163:A:MET:HE1	1:164:A:GLY:HA2	3	0.3
(1,519)	1:163:A:MET:HE2	1:164:A:GLY:HA2	3	0.3
(1,519)	1:163:A:MET:HE3	1:164:A:GLY:HA2	3	0.3
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB1	5	0.29
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB2	5	0.29
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB3	5	0.29
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB1	5	0.29
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB2	5	0.29
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB3	5	0.29
(1,1644)	1:108:A:LEU:HD11	1:118:A:GLU:HG2	6	0.29
(1,1644)	1:108:A:LEU:HD11	1:118:A:GLU:HG3	6	0.29
(1,1644)	1:108:A:LEU:HD12	1:118:A:GLU:HG2	6	0.29
(1,1644)	1:108:A:LEU:HD12	1:118:A:GLU:HG3	6	0.29
(1,1644)	1:108:A:LEU:HD13	1:118:A:GLU:HG2	6	0.29
(1,1644)	1:108:A:LEU:HD13	1:118:A:GLU:HG3	6	0.29
(1,1644)	1:108:A:LEU:HD21	1:118:A:GLU:HG2	6	0.29
(1,1644)	1:108:A:LEU:HD21	1:118:A:GLU:HG3	6	0.29
(1,1644)	1:108:A:LEU:HD22	1:118:A:GLU:HG2	6	0.29
(1,1644)	1:108:A:LEU:HD22	1:118:A:GLU:HG3	6	0.29
(1,1644)	1:108:A:LEU:HD23	1:118:A:GLU:HG2	6	0.29
(1,1644)	1:108:A:LEU:HD23	1:118:A:GLU:HG3	6	0.29
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD11	9	0.29
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD12	9	0.29

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD13	9	0.29
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD21	9	0.29
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD22	9	0.29
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD23	9	0.29
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD11	9	0.29
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD12	9	0.29
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD13	9	0.29
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD21	9	0.29
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD22	9	0.29
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD23	9	0.29
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	9	0.29
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	9	0.29
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	9	0.29
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	9	0.29
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	9	0.29
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	9	0.29
(1,795)	1:83:A:GLU:H	1:84:A:LEU:HG	14	0.29
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD21	13	0.28
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD22	13	0.28
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD21	13	0.28
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD22	13	0.28
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD21	13	0.28
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD22	13	0.28
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD21	13	0.28
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD22	13	0.28
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD21	13	0.28
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD22	13	0.28
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD21	13	0.28
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD22	13	0.28
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG2	7	0.28
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG3	7	0.28
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	5	0.28
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	5	0.28
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	5	0.28
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	5	0.28
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	5	0.28
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	5	0.28
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	3	0.27
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	3	0.27
(1,1682)	1:121:A:THR:H	1:124:A:GLU:HG2	3	0.27
(1,1682)	1:121:A:THR:H	1:124:A:GLU:HG3	3	0.27
(1,1134)	1:135:A:LYS:HG2	1:150:A:GLN:H	5	0.27

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,210)	1:59:A:ALA:HA	1:66:A:THR:HA	12	0.27
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG21	4	0.27
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG22	4	0.27
(1,104)	1:44:A:LEU:HG	1:46:A:VAL:HG23	4	0.27
(1,2)	1:59:A:ALA:HB1	1:66:A:THR:HA	6	0.27
(1,2)	1:59:A:ALA:HB2	1:66:A:THR:HA	6	0.27
(1,2)	1:59:A:ALA:HB3	1:66:A:THR:HA	6	0.27
(1,1753)	1:134:A:TYR:HE1	1:149:A:GLN:HG2	13	0.26
(1,1753)	1:134:A:TYR:HE1	1:149:A:GLN:HG3	13	0.26
(1,1753)	1:134:A:TYR:HE2	1:149:A:GLN:HG2	13	0.26
(1,1753)	1:134:A:TYR:HE2	1:149:A:GLN:HG3	13	0.26
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG21	6	0.26
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG22	6	0.26
(1,1443)	1:53:A:GLY:HA2	1:54:A:ILE:HG23	6	0.26
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG21	6	0.26
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG22	6	0.26
(1,1443)	1:53:A:GLY:HA3	1:54:A:ILE:HG23	6	0.26
(1,555)	1:134:A:TYR:HE1	1:149:A:GLN:HG2	13	0.26
(1,555)	1:134:A:TYR:HE2	1:149:A:GLN:HG2	13	0.26
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD21	4	0.26
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD22	4	0.26
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD23	4	0.26
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	11	0.25
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	11	0.25
(2,186)	1:81:A:LEU:HD11	1:130:A:LEU:H	13	0.25
(2,186)	1:81:A:LEU:HD12	1:130:A:LEU:H	13	0.25
(2,186)	1:81:A:LEU:HD13	1:130:A:LEU:H	13	0.25
(2,186)	1:81:A:LEU:HD21	1:130:A:LEU:H	13	0.25
(2,186)	1:81:A:LEU:HD22	1:130:A:LEU:H	13	0.25
(2,186)	1:81:A:LEU:HD23	1:130:A:LEU:H	13	0.25
(2,76)	1:135:A:LYS:HE2	1:150:A:GLN:HB2	1	0.25
(2,76)	1:135:A:LYS:HE2	1:150:A:GLN:HB3	1	0.25
(2,76)	1:135:A:LYS:HE3	1:150:A:GLN:HB2	1	0.25
(2,76)	1:135:A:LYS:HE3	1:150:A:GLN:HB3	1	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD11	5	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD12	5	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD13	5	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD21	5	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD22	5	0.25
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD23	5	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD11	5	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD12	5	0.25

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD13	5	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD21	5	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD22	5	0.25
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD23	5	0.25
(1,941)	1:40:A:ALA:HA	1:107:A:ASP:H	13	0.25
(1,78)	1:137:A:VAL:HG21	1:147:A:LEU:HG	14	0.25
(1,78)	1:137:A:VAL:HG22	1:147:A:LEU:HG	14	0.25
(1,78)	1:137:A:VAL:HG23	1:147:A:LEU:HG	14	0.25
(2,136)	1:2:A:LEU:HG	1:166:A:LEU:H	9	0.24
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD21	4	0.24
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD22	4	0.24
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD23	4	0.24
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD21	13	0.24
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD22	13	0.24
(2,20)	1:85:A:ASP:H	1:132:A:LEU:HD23	13	0.24
(1,981)	1:120:A:ARG:H	1:120:A:ARG:HD3	3	0.24
(1,622)	1:46:A:VAL:HA	1:61:A:LEU:HB2	14	0.24
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG21	8	0.24
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG22	8	0.24
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG23	8	0.24
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD21	11	0.23
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD22	11	0.23
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD21	11	0.23
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD22	11	0.23
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD21	11	0.23
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD22	11	0.23
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD21	11	0.23
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD22	11	0.23
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD21	11	0.23
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD22	11	0.23
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD21	11	0.23
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD22	11	0.23
(2,198)	1:103:A:ARG:HD2	1:122:A:LEU:H	6	0.23
(2,198)	1:103:A:ARG:HD3	1:122:A:LEU:H	6	0.23
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	9	0.23
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	9	0.23
(2,113)	1:90:A:ASN:H	1:132:A:LEU:HG	13	0.23
(1,1650)	1:114:A:SER:HB2	1:119:A:LEU:HG	4	0.23
(1,1650)	1:114:A:SER:HB3	1:119:A:LEU:HG	4	0.23
(1,1414)	1:45:A:LEU:HD11	1:163:A:MET:HE1	7	0.23
(1,1414)	1:45:A:LEU:HD11	1:163:A:MET:HE2	7	0.23
(1,1414)	1:45:A:LEU:HD11	1:163:A:MET:HE3	7	0.23

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1414)	1:45:A:LEU:HD12	1:163:A:MET:HE1	7	0.23
(1,1414)	1:45:A:LEU:HD12	1:163:A:MET:HE2	7	0.23
(1,1414)	1:45:A:LEU:HD12	1:163:A:MET:HE3	7	0.23
(1,1414)	1:45:A:LEU:HD13	1:163:A:MET:HE1	7	0.23
(1,1414)	1:45:A:LEU:HD13	1:163:A:MET:HE2	7	0.23
(1,1414)	1:45:A:LEU:HD13	1:163:A:MET:HE3	7	0.23
(1,1414)	1:45:A:LEU:HD21	1:163:A:MET:HE1	7	0.23
(1,1414)	1:45:A:LEU:HD21	1:163:A:MET:HE2	7	0.23
(1,1414)	1:45:A:LEU:HD21	1:163:A:MET:HE3	7	0.23
(1,1414)	1:45:A:LEU:HD22	1:163:A:MET:HE1	7	0.23
(1,1414)	1:45:A:LEU:HD22	1:163:A:MET:HE2	7	0.23
(1,1414)	1:45:A:LEU:HD22	1:163:A:MET:HE3	7	0.23
(1,1414)	1:45:A:LEU:HD23	1:163:A:MET:HE1	7	0.23
(1,1414)	1:45:A:LEU:HD23	1:163:A:MET:HE2	7	0.23
(1,1414)	1:45:A:LEU:HD23	1:163:A:MET:HE3	7	0.23
(1,112)	1:46:A:VAL:HG11	1:55:A:ILE:HB	7	0.23
(1,112)	1:46:A:VAL:HG12	1:55:A:ILE:HB	7	0.23
(1,112)	1:46:A:VAL:HG13	1:55:A:ILE:HB	7	0.23
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	8	0.22
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	8	0.22
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD11	3	0.22
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD12	3	0.22
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD13	3	0.22
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD21	3	0.22
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD22	3	0.22
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD23	3	0.22
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD11	3	0.22
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD12	3	0.22
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD13	3	0.22
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD21	3	0.22
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD22	3	0.22
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD23	3	0.22
(1,218)	1:68:A:LEU:HB3	1:83:A:GLU:HG2	6	0.22
(1,1643)	1:108:A:LEU:HD11	1:118:A:GLU:HA	1	0.21
(1,1643)	1:108:A:LEU:HD12	1:118:A:GLU:HA	1	0.21
(1,1643)	1:108:A:LEU:HD13	1:118:A:GLU:HA	1	0.21
(1,1643)	1:108:A:LEU:HD21	1:118:A:GLU:HA	1	0.21
(1,1643)	1:108:A:LEU:HD22	1:118:A:GLU:HA	1	0.21
(1,1643)	1:108:A:LEU:HD23	1:118:A:GLU:HA	1	0.21
(1,1480)	1:68:A:LEU:HD11	1:75:A:LYS:H	6	0.21
(1,1480)	1:68:A:LEU:HD12	1:75:A:LYS:H	6	0.21
(1,1480)	1:68:A:LEU:HD13	1:75:A:LYS:H	6	0.21

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1480)	1:68:A:LEU:HD21	1:75:A:LYS:H	6	0.21
(1,1480)	1:68:A:LEU:HD22	1:75:A:LYS:H	6	0.21
(1,1480)	1:68:A:LEU:HD23	1:75:A:LYS:H	6	0.21
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD21	13	0.21
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD22	13	0.21
(1,338)	1:82:A:ALA:HA	1:132:A:LEU:HD23	13	0.21
(1,218)	1:68:A:LEU:HB3	1:83:A:GLU:HG2	14	0.21
(1,123)	1:47:A:ARG:H	1:49:A:LEU:HD11	5	0.21
(1,123)	1:47:A:ARG:H	1:49:A:LEU:HD12	5	0.21
(1,123)	1:47:A:ARG:H	1:49:A:LEU:HD13	5	0.21
(1,93)	1:40:A:ALA:HB1	1:108:A:LEU:HA	7	0.21
(1,93)	1:40:A:ALA:HB2	1:108:A:LEU:HA	7	0.21
(1,93)	1:40:A:ALA:HB3	1:108:A:LEU:HA	7	0.21
(1,93)	1:40:A:ALA:HB1	1:108:A:LEU:HA	8	0.21
(1,93)	1:40:A:ALA:HB2	1:108:A:LEU:HA	8	0.21
(1,93)	1:40:A:ALA:HB3	1:108:A:LEU:HA	8	0.21
(2,65)	1:0:A:ASP:HA	1:61:A:LEU:HD11	1	0.2
(2,65)	1:0:A:ASP:HA	1:61:A:LEU:HD12	1	0.2
(2,65)	1:0:A:ASP:HA	1:61:A:LEU:HD13	1	0.2
(2,65)	1:0:A:ASP:HA	1:61:A:LEU:HD21	1	0.2
(2,65)	1:0:A:ASP:HA	1:61:A:LEU:HD22	1	0.2
(2,65)	1:0:A:ASP:HA	1:61:A:LEU:HD23	1	0.2
(1,1420)	1:47:A:ARG:HB2	1:49:A:LEU:H	5	0.2
(1,1420)	1:47:A:ARG:HB3	1:49:A:LEU:H	5	0.2
(1,1354)	1:10:A:GLY:H	1:162:A:LYS:HG2	14	0.2
(1,1354)	1:10:A:GLY:H	1:162:A:LYS:HG3	14	0.2
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD21	2	0.19
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD22	2	0.19
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD21	2	0.19
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD22	2	0.19
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD21	2	0.19
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD22	2	0.19
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD21	2	0.19
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD22	2	0.19
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD21	2	0.19
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD22	2	0.19
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD21	2	0.19
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD22	2	0.19
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD11	1	0.19
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD12	1	0.19
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD13	1	0.19
(1,1566)	1:90:A:ASN:HB3	1:91:A:ILE:HG12	13	0.19

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1566)	1:90:A:ASN:HB3	1:91:A:ILE:HG13	13	0.19
(1,1101)	1:139:A:LYS:H	1:147:A:LEU:HG	14	0.19
(1,885)	1:47:A:ARG:H	1:49:A:LEU:HD21	7	0.19
(1,885)	1:47:A:ARG:H	1:49:A:LEU:HD22	7	0.19
(1,885)	1:47:A:ARG:H	1:49:A:LEU:HD23	7	0.19
(1,519)	1:163:A:MET:HE1	1:164:A:GLY:HA2	12	0.19
(1,519)	1:163:A:MET:HE2	1:164:A:GLY:HA2	12	0.19
(1,519)	1:163:A:MET:HE3	1:164:A:GLY:HA2	12	0.19
(1,313)	1:81:A:LEU:HD11	1:92:A:ILE:HD11	1	0.19
(1,313)	1:81:A:LEU:HD11	1:92:A:ILE:HD12	1	0.19
(1,313)	1:81:A:LEU:HD11	1:92:A:ILE:HD13	1	0.19
(1,313)	1:81:A:LEU:HD12	1:92:A:ILE:HD11	1	0.19
(1,313)	1:81:A:LEU:HD12	1:92:A:ILE:HD12	1	0.19
(1,313)	1:81:A:LEU:HD12	1:92:A:ILE:HD13	1	0.19
(1,313)	1:81:A:LEU:HD13	1:92:A:ILE:HD11	1	0.19
(1,313)	1:81:A:LEU:HD13	1:92:A:ILE:HD12	1	0.19
(1,313)	1:81:A:LEU:HD13	1:92:A:ILE:HD13	1	0.19
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	7	0.18
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	7	0.18
(2,75)	1:135:A:LYS:HE2	1:137:A:VAL:HB	1	0.18
(2,75)	1:135:A:LYS:HE3	1:137:A:VAL:HB	1	0.18
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD11	11	0.18
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD12	11	0.18
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD13	11	0.18
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD21	11	0.18
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD22	11	0.18
(1,1636)	1:106:A:GLN:HE21	1:108:A:LEU:HD23	11	0.18
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD11	11	0.18
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD12	11	0.18
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD13	11	0.18
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD21	11	0.18
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD22	11	0.18
(1,1636)	1:106:A:GLN:HE22	1:108:A:LEU:HD23	11	0.18
(1,1420)	1:47:A:ARG:HB2	1:49:A:LEU:H	3	0.18
(1,1420)	1:47:A:ARG:HB3	1:49:A:LEU:H	3	0.18
(1,859)	1:93:A:ASN:HD21	1:94:A:ALA:H	4	0.18
(1,736)	1:61:A:LEU:H	1:61:A:LEU:HG	8	0.18
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD21	13	0.18
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD22	13	0.18
(1,337)	1:82:A:ALA:HB1	1:132:A:LEU:HD23	13	0.18
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD21	13	0.18
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD22	13	0.18

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,337)	1:82:A:ALA:HB2	1:132:A:LEU:HD23	13	0.18
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD21	13	0.18
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD22	13	0.18
(1,337)	1:82:A:ALA:HB3	1:132:A:LEU:HD23	13	0.18
(2,212)	1:131:A:ASP:HB2	1:154:A:THR:H	12	0.17
(2,212)	1:131:A:ASP:HB3	1:154:A:THR:H	12	0.17
(2,136)	1:2:A:LEU:HG	1:166:A:LEU:H	1	0.17
(2,50)	1:152:A:SER:HA	1:160:A:THR:HB	14	0.17
(1,1500)	1:80:A:ALA:HA	1:83:A:GLU:HG2	8	0.17
(1,1500)	1:80:A:ALA:HA	1:83:A:GLU:HG3	8	0.17
(1,622)	1:46:A:VAL:HA	1:61:A:LEU:HB2	7	0.17
(1,272)	1:94:A:ALA:HA	1:100:A:GLN:HG2	5	0.17
(1,272)	1:94:A:ALA:HA	1:100:A:GLN:HG3	5	0.17
(2,75)	1:135:A:LYS:HE2	1:137:A:VAL:HB	10	0.16
(2,75)	1:135:A:LYS:HE3	1:137:A:VAL:HB	10	0.16
(1,1360)	1:11:A:ILE:HG12	1:163:A:MET:HB3	11	0.16
(1,1360)	1:11:A:ILE:HG13	1:163:A:MET:HB3	11	0.16
(2,217)	1:141:A:LEU:HD11	1:145:A:ASN:H	5	0.15
(2,217)	1:141:A:LEU:HD12	1:145:A:ASN:H	5	0.15
(2,217)	1:141:A:LEU:HD13	1:145:A:ASN:H	5	0.15
(2,217)	1:141:A:LEU:HD21	1:145:A:ASN:H	5	0.15
(2,217)	1:141:A:LEU:HD22	1:145:A:ASN:H	5	0.15
(2,217)	1:141:A:LEU:HD23	1:145:A:ASN:H	5	0.15
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	2	0.15
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	2	0.15
(2,186)	1:81:A:LEU:HD11	1:130:A:LEU:H	12	0.15
(2,186)	1:81:A:LEU:HD12	1:130:A:LEU:H	12	0.15
(2,186)	1:81:A:LEU:HD13	1:130:A:LEU:H	12	0.15
(2,186)	1:81:A:LEU:HD21	1:130:A:LEU:H	12	0.15
(2,186)	1:81:A:LEU:HD22	1:130:A:LEU:H	12	0.15
(2,186)	1:81:A:LEU:HD23	1:130:A:LEU:H	12	0.15
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB1	7	0.15
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB2	7	0.15
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB3	7	0.15
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	7	0.15
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	7	0.15
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	7	0.15
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG2	4	0.15
(1,1471)	1:68:A:LEU:HG	1:83:A:GLU:HG3	4	0.15
(1,1422)	1:47:A:ARG:HB2	1:49:A:LEU:HD11	3	0.15
(1,1422)	1:47:A:ARG:HB2	1:49:A:LEU:HD12	3	0.15
(1,1422)	1:47:A:ARG:HB2	1:49:A:LEU:HD13	3	0.15

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1422)	1:47:A:ARG:HB2	1:49:A:LEU:HD21	3	0.15
(1,1422)	1:47:A:ARG:HB2	1:49:A:LEU:HD22	3	0.15
(1,1422)	1:47:A:ARG:HB2	1:49:A:LEU:HD23	3	0.15
(1,1422)	1:47:A:ARG:HB3	1:49:A:LEU:HD11	3	0.15
(1,1422)	1:47:A:ARG:HB3	1:49:A:LEU:HD12	3	0.15
(1,1422)	1:47:A:ARG:HB3	1:49:A:LEU:HD13	3	0.15
(1,1422)	1:47:A:ARG:HB3	1:49:A:LEU:HD21	3	0.15
(1,1422)	1:47:A:ARG:HB3	1:49:A:LEU:HD22	3	0.15
(1,1422)	1:47:A:ARG:HB3	1:49:A:LEU:HD23	3	0.15
(1,1419)	1:47:A:ARG:H	1:49:A:LEU:HD11	5	0.15
(1,1419)	1:47:A:ARG:H	1:49:A:LEU:HD12	5	0.15
(1,1419)	1:47:A:ARG:H	1:49:A:LEU:HD13	5	0.15
(1,1419)	1:47:A:ARG:H	1:49:A:LEU:HD21	5	0.15
(1,1419)	1:47:A:ARG:H	1:49:A:LEU:HD22	5	0.15
(1,1419)	1:47:A:ARG:H	1:49:A:LEU:HD23	5	0.15
(1,736)	1:61:A:LEU:H	1:61:A:LEU:HG	6	0.15
(2,198)	1:103:A:ARG:HD2	1:122:A:LEU:H	3	0.14
(2,198)	1:103:A:ARG:HD3	1:122:A:LEU:H	3	0.14
(2,137)	1:9:A:ASP:H	1:162:A:LYS:H	10	0.14
(2,68)	1:0:A:ASP:HA	1:166:A:LEU:HB2	13	0.14
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB1	1	0.14
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB2	1	0.14
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB3	1	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	1	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	1	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	1	0.14
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB1	11	0.14
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB2	11	0.14
(1,1556)	1:87:A:ASN:HD21	1:95:A:ALA:HB3	11	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB1	11	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB2	11	0.14
(1,1556)	1:87:A:ASN:HD22	1:95:A:ALA:HB3	11	0.14
(1,1420)	1:47:A:ARG:HB2	1:49:A:LEU:H	11	0.14
(1,1420)	1:47:A:ARG:HB3	1:49:A:LEU:H	11	0.14
(1,1360)	1:11:A:ILE:HG12	1:163:A:MET:HB3	9	0.14
(1,1360)	1:11:A:ILE:HG13	1:163:A:MET:HB3	9	0.14
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE1	7	0.14
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE2	7	0.14
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE3	7	0.14
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE1	7	0.14
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE2	7	0.14
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE3	7	0.14

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE1	7	0.14
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE2	7	0.14
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE3	7	0.14
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE1	7	0.14
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE2	7	0.14
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE3	7	0.14
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE1	7	0.14
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE2	7	0.14
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE3	7	0.14
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE1	7	0.14
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE2	7	0.14
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE3	7	0.14
(1,623)	1:46:A:VAL:HB	1:55:A:ILE:HG21	7	0.14
(1,623)	1:46:A:VAL:HB	1:55:A:ILE:HG22	7	0.14
(1,623)	1:46:A:VAL:HB	1:55:A:ILE:HG23	7	0.14
(2,191)	1:87:A:ASN:HD21	1:96:A:ASP:H	10	0.13
(2,191)	1:87:A:ASN:HD22	1:96:A:ASP:H	10	0.13
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD11	14	0.13
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD12	14	0.13
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD13	14	0.13
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD21	14	0.13
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD22	14	0.13
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD23	14	0.13
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD21	6	0.13
(1,1642)	1:108:A:LEU:HD11	1:117:A:ASN:HD22	6	0.13
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD21	6	0.13
(1,1642)	1:108:A:LEU:HD12	1:117:A:ASN:HD22	6	0.13
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD21	6	0.13
(1,1642)	1:108:A:LEU:HD13	1:117:A:ASN:HD22	6	0.13
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD21	6	0.13
(1,1642)	1:108:A:LEU:HD21	1:117:A:ASN:HD22	6	0.13
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD21	6	0.13
(1,1642)	1:108:A:LEU:HD22	1:117:A:ASN:HD22	6	0.13
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD21	6	0.13
(1,1642)	1:108:A:LEU:HD23	1:117:A:ASN:HD22	6	0.13
(1,1566)	1:90:A:ASN:HB3	1:91:A:ILE:HG12	1	0.13
(1,1566)	1:90:A:ASN:HB3	1:91:A:ILE:HG13	1	0.13
(1,1470)	1:68:A:LEU:HB3	1:83:A:GLU:HB2	12	0.13
(1,1470)	1:68:A:LEU:HB3	1:83:A:GLU:HB3	12	0.13
(1,1283)	1:1:A:ALA:HB1	1:168:A:LEU:HD11	14	0.13
(1,1283)	1:1:A:ALA:HB1	1:168:A:LEU:HD12	14	0.13
(1,1283)	1:1:A:ALA:HB1	1:168:A:LEU:HD13	14	0.13

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1283)	1:1:A:ALA:HB1	1:168:A:LEU:HD21	14	0.13
(1,1283)	1:1:A:ALA:HB1	1:168:A:LEU:HD22	14	0.13
(1,1283)	1:1:A:ALA:HB1	1:168:A:LEU:HD23	14	0.13
(1,1283)	1:1:A:ALA:HB2	1:168:A:LEU:HD11	14	0.13
(1,1283)	1:1:A:ALA:HB2	1:168:A:LEU:HD12	14	0.13
(1,1283)	1:1:A:ALA:HB2	1:168:A:LEU:HD13	14	0.13
(1,1283)	1:1:A:ALA:HB2	1:168:A:LEU:HD21	14	0.13
(1,1283)	1:1:A:ALA:HB2	1:168:A:LEU:HD22	14	0.13
(1,1283)	1:1:A:ALA:HB2	1:168:A:LEU:HD23	14	0.13
(1,1283)	1:1:A:ALA:HB3	1:168:A:LEU:HD11	14	0.13
(1,1283)	1:1:A:ALA:HB3	1:168:A:LEU:HD12	14	0.13
(1,1283)	1:1:A:ALA:HB3	1:168:A:LEU:HD13	14	0.13
(1,1283)	1:1:A:ALA:HB3	1:168:A:LEU:HD21	14	0.13
(1,1283)	1:1:A:ALA:HB3	1:168:A:LEU:HD22	14	0.13
(1,1283)	1:1:A:ALA:HB3	1:168:A:LEU:HD23	14	0.13
(1,1191)	1:159:A:THR:HB	1:160:A:THR:H	1	0.13
(1,471)	1:134:A:TYR:HE1	1:149:A:GLN:HG3	3	0.13
(1,471)	1:134:A:TYR:HE2	1:149:A:GLN:HG3	3	0.13
(1,62)	1:11:A:ILE:HD11	1:162:A:LYS:HA	2	0.13
(1,62)	1:11:A:ILE:HD12	1:162:A:LYS:HA	2	0.13
(1,62)	1:11:A:ILE:HD13	1:162:A:LYS:HA	2	0.13
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD21	9	0.12
(2,204)	1:108:A:LEU:HD11	1:109:A:ASN:HD22	9	0.12
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD21	9	0.12
(2,204)	1:108:A:LEU:HD12	1:109:A:ASN:HD22	9	0.12
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD21	9	0.12
(2,204)	1:108:A:LEU:HD13	1:109:A:ASN:HD22	9	0.12
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD21	9	0.12
(2,204)	1:108:A:LEU:HD21	1:109:A:ASN:HD22	9	0.12
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD21	9	0.12
(2,204)	1:108:A:LEU:HD22	1:109:A:ASN:HD22	9	0.12
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD21	9	0.12
(2,204)	1:108:A:LEU:HD23	1:109:A:ASN:HD22	9	0.12
(2,187)	1:82:A:ALA:H	1:83:A:GLU:HG2	2	0.12
(2,187)	1:82:A:ALA:H	1:83:A:GLU:HG3	2	0.12
(2,187)	1:82:A:ALA:H	1:83:A:GLU:HG2	9	0.12
(2,187)	1:82:A:ALA:H	1:83:A:GLU:HG3	9	0.12
(2,68)	1:0:A:ASP:HA	1:166:A:LEU:HB2	8	0.12
(1,1561)	1:90:A:ASN:H	1:132:A:LEU:HD11	13	0.12
(1,1561)	1:90:A:ASN:H	1:132:A:LEU:HD12	13	0.12
(1,1561)	1:90:A:ASN:H	1:132:A:LEU:HD13	13	0.12
(1,1561)	1:90:A:ASN:H	1:132:A:LEU:HD21	13	0.12

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1561)	1:90:A:ASN:H	1:132:A:LEU:HD22	13	0.12
(1,1561)	1:90:A:ASN:H	1:132:A:LEU:HD23	13	0.12
(1,622)	1:46:A:VAL:HA	1:61:A:LEU:HB2	10	0.12
(1,521)	1:163:A:MET:HE1	1:164:A:GLY:H	7	0.12
(1,521)	1:163:A:MET:HE2	1:164:A:GLY:H	7	0.12
(1,521)	1:163:A:MET:HE3	1:164:A:GLY:H	7	0.12
(1,323)	1:84:A:LEU:HB3	1:92:A:ILE:HG21	4	0.12
(1,323)	1:84:A:LEU:HB3	1:92:A:ILE:HG22	4	0.12
(1,323)	1:84:A:LEU:HB3	1:92:A:ILE:HG23	4	0.12
(1,272)	1:94:A:ALA:HA	1:100:A:GLN:HG2	13	0.12
(1,272)	1:94:A:ALA:HA	1:100:A:GLN:HG3	13	0.12
(2,76)	1:135:A:LYS:HE2	1:150:A:GLN:HB2	10	0.11
(2,76)	1:135:A:LYS:HE2	1:150:A:GLN:HB3	10	0.11
(2,76)	1:135:A:LYS:HE3	1:150:A:GLN:HB2	10	0.11
(2,76)	1:135:A:LYS:HE3	1:150:A:GLN:HB3	10	0.11
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB1	12	0.11
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB2	12	0.11
(2,61)	1:134:A:TYR:HE1	1:148:A:ALA:HB3	12	0.11
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB1	12	0.11
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB2	12	0.11
(2,61)	1:134:A:TYR:HE2	1:148:A:ALA:HB3	12	0.11
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD11	11	0.11
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD12	11	0.11
(2,9)	1:90:A:ASN:HB3	1:91:A:ILE:HD13	11	0.11
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE1	8	0.11
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE2	8	0.11
(1,1330)	1:4:A:LEU:HD11	1:163:A:MET:HE3	8	0.11
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE1	8	0.11
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE2	8	0.11
(1,1330)	1:4:A:LEU:HD12	1:163:A:MET:HE3	8	0.11
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE1	8	0.11
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE2	8	0.11
(1,1330)	1:4:A:LEU:HD13	1:163:A:MET:HE3	8	0.11
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE1	8	0.11
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE2	8	0.11
(1,1330)	1:4:A:LEU:HD21	1:163:A:MET:HE3	8	0.11
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE1	8	0.11
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE2	8	0.11
(1,1330)	1:4:A:LEU:HD22	1:163:A:MET:HE3	8	0.11
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE1	8	0.11
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE2	8	0.11
(1,1330)	1:4:A:LEU:HD23	1:163:A:MET:HE3	8	0.11

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,736)	1:61:A:LEU:H	1:61:A:LEU:HG	3	0.11
(1,218)	1:68:A:LEU:HB3	1:83:A:GLU:HG2	3	0.11
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG21	5	0.11
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG22	5	0.11
(1,79)	1:135:A:LYS:HB3	1:137:A:VAL:HG23	5	0.11
(2,174)	1:46:A:VAL:HA	1:49:A:LEU:HD11	5	0.1
(2,174)	1:46:A:VAL:HA	1:49:A:LEU:HD12	5	0.1
(2,174)	1:46:A:VAL:HA	1:49:A:LEU:HD13	5	0.1
(2,174)	1:46:A:VAL:HA	1:49:A:LEU:HD21	5	0.1
(2,174)	1:46:A:VAL:HA	1:49:A:LEU:HD22	5	0.1
(2,174)	1:46:A:VAL:HA	1:49:A:LEU:HD23	5	0.1
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD11	5	0.1
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD12	5	0.1
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD13	5	0.1
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD21	5	0.1
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD22	5	0.1
(1,1828)	1:166:A:LEU:HB2	1:168:A:LEU:HD23	5	0.1
(1,266)	1:99:A:PHE:HA	1:102:A:LEU:HD21	12	0.1
(1,266)	1:99:A:PHE:HA	1:102:A:LEU:HD22	12	0.1
(1,266)	1:99:A:PHE:HA	1:102:A:LEU:HD23	12	0.1
(1,210)	1:59:A:ALA:HA	1:66:A:THR:HA	8	0.1

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

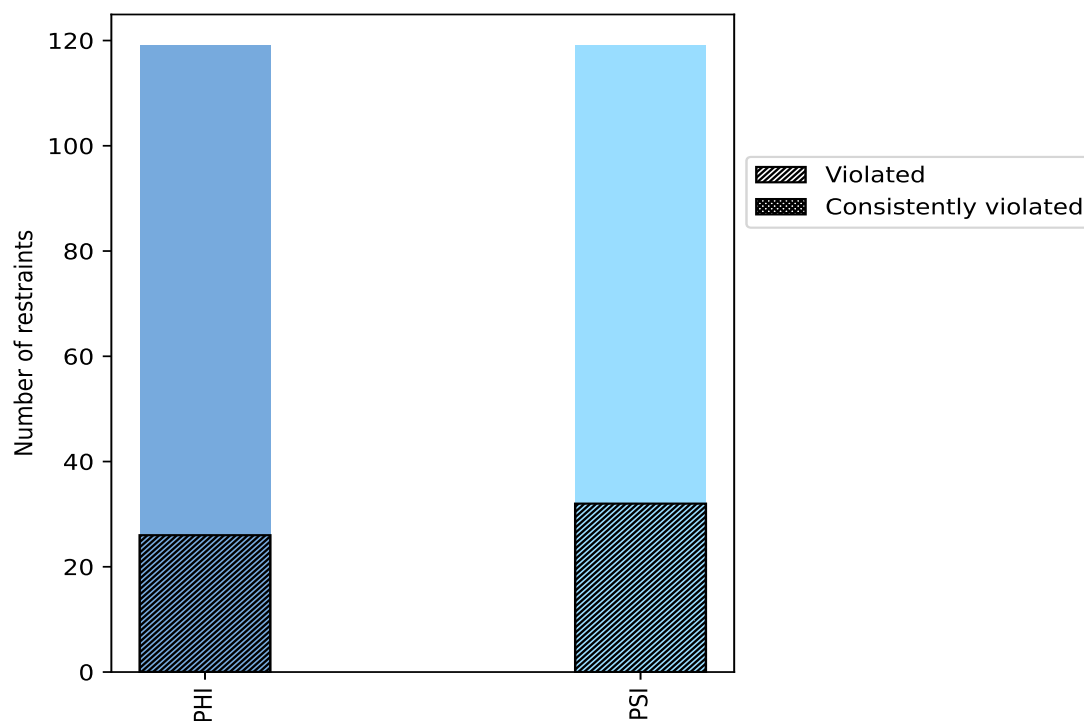
10.1 Summary of dihedral-angle violations [i](#)

The following table provides the summary of dihedral-angle violations in different dihedral-angle types. Violations less than 1° are not included in the calculation.

Angle type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
PHI	119	50.0	26	21.8	10.9	0	0.0	0.0
PSI	119	50.0	32	26.9	13.4	0	0.0	0.0
Total	238	100.0	58	24.4	24.4	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](#)



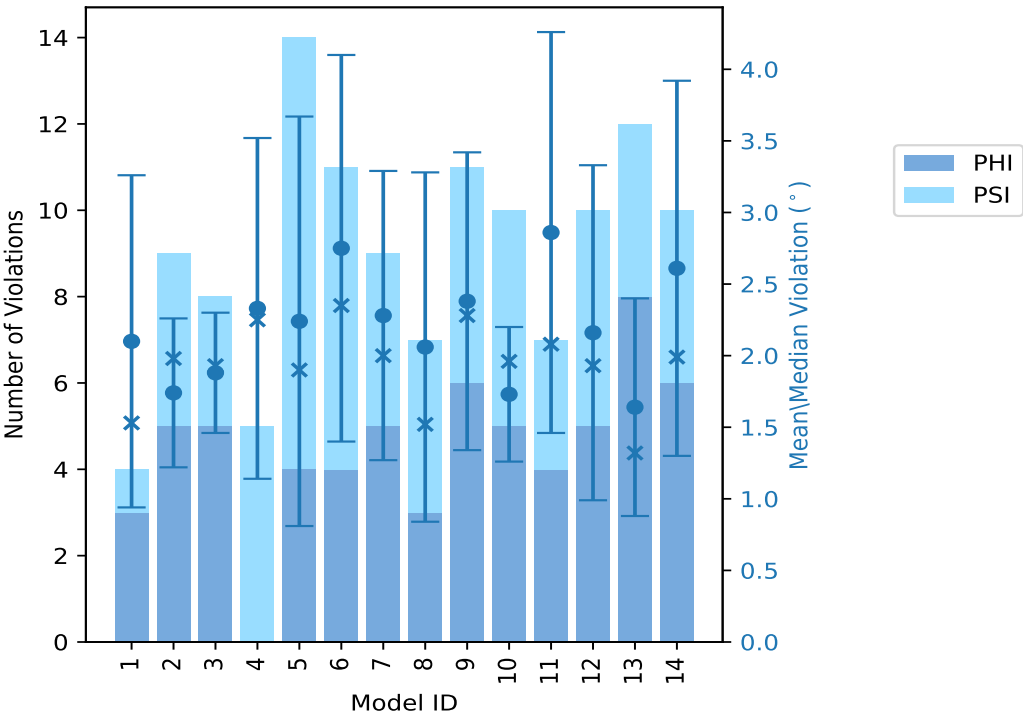
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

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 (°)
	PHI	PSI	Total				
1	3	1	4	2.1	4.1	1.16	1.53
2	5	4	9	1.74	2.45	0.52	1.98
3	5	3	8	1.88	2.55	0.42	1.93
4	0	5	5	2.33	4.16	1.19	2.25
5	4	10	14	2.24	6.8	1.43	1.9
6	4	7	11	2.75	5.15	1.35	2.35
7	5	4	9	2.28	4.33	1.01	2.0
8	3	4	7	2.06	4.71	1.22	1.52
9	6	5	11	2.38	4.03	1.04	2.28
10	5	5	10	1.73	2.37	0.47	1.96
11	4	3	7	2.86	5.28	1.4	2.08
12	5	5	10	2.16	5.33	1.17	1.93
13	8	4	12	1.64	3.88	0.76	1.32
14	6	4	10	2.61	5.22	1.31	1.99

10.2.1 Bar graph : Dihedral 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

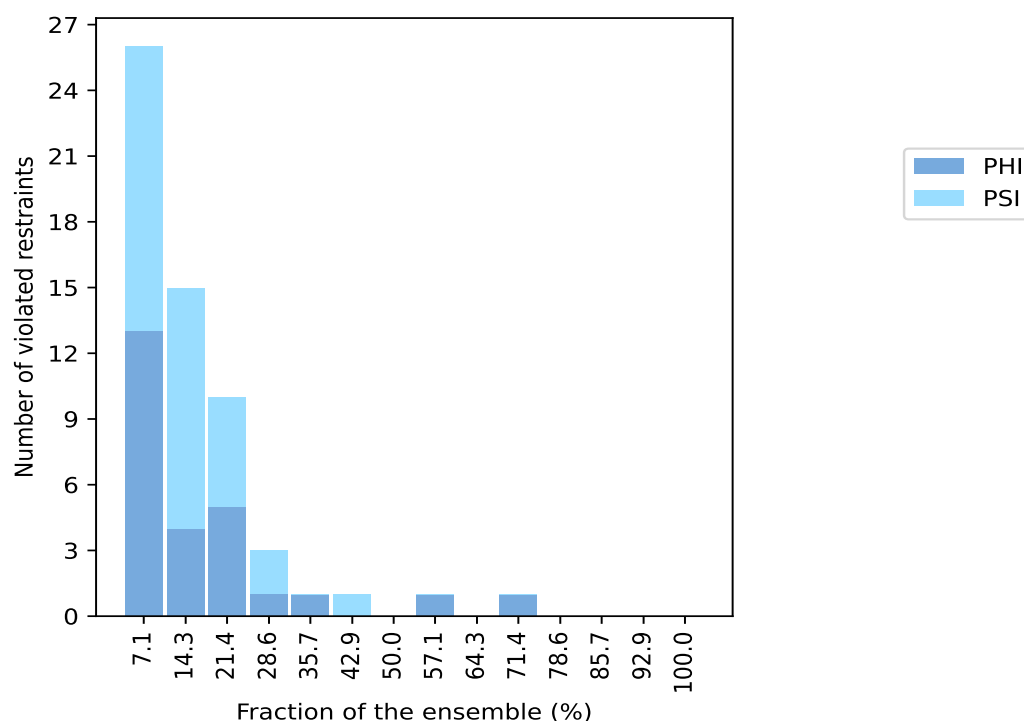
10.3 Dihedral-angle violation statistics for the ensemble

Violation analysis may find that some restraints are violated in very few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of ensemble.

Number of violated restraints			Fraction of the ensemble	
PHI	PSI	Total	Count ¹	%
13	13	26	1	7.1
4	11	15	2	14.3
5	5	10	3	21.4
1	2	3	4	28.6
1	0	1	5	35.7
0	1	1	6	42.9
0	0	0	7	50.0
1	0	1	8	57.1
0	0	0	9	64.3
1	0	1	10	71.4
0	0	0	11	78.6
0	0	0	12	85.7
0	0	0	13	92.9
0	0	0	14	100.0

¹ Number of models with violations

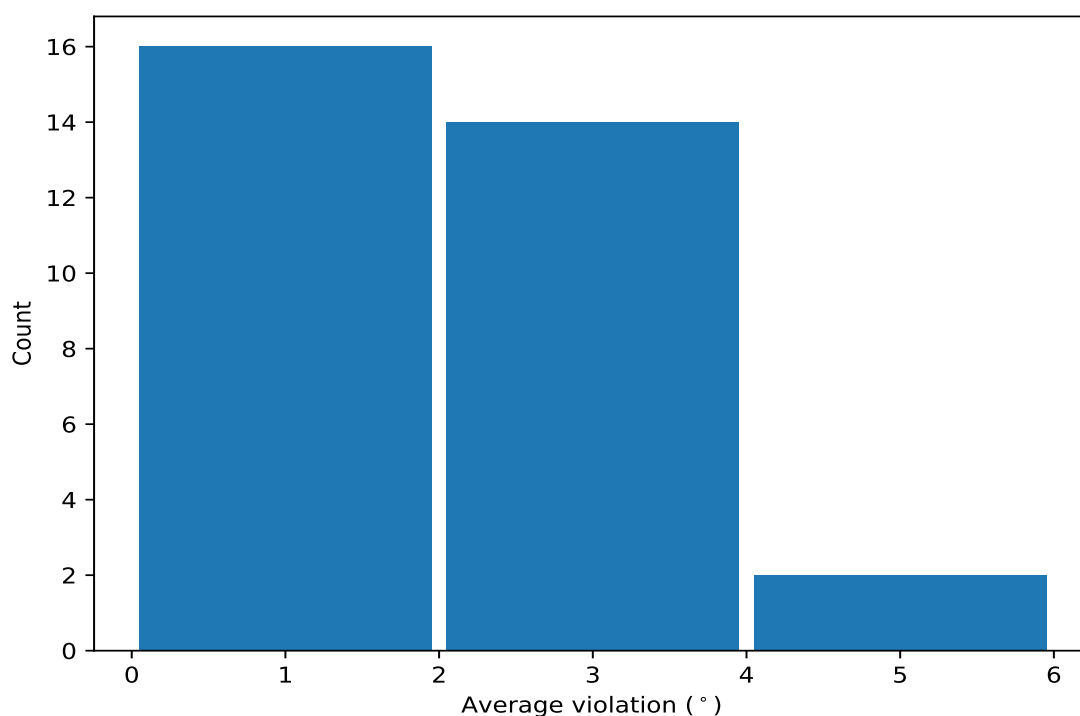
10.3.1 Bar graph : Dihedral-angle Violation statistics for the ensemble [i](#)



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
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	10	3.75	1.63	3.84
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	8	1.9	0.51	1.74
(1,134)	1:106:A:GLN:N	1:106:A:GLN:CA	1:106:A:GLN:C	1:107:A:ASP:N	6	1.88	0.44	1.98
(1,11)	1:5:A:ASP:C	1:6:A:LEU:N	1:6:A:LEU:CA	1:6:A:LEU:C	5	2.13	0.97	1.72
(1,226)	1:165:A:ASP:N	1:165:A:ASP:CA	1:165:A:ASP:C	1:166:A:LEU:N	4	2.08	0.73	2.06
(1,45)	1:48:A:ASP:C	1:49:A:LEU:N	1:49:A:LEU:CA	1:49:A:LEU:C	4	1.64	0.53	1.6
(1,32)	1:40:A:ALA:N	1:40:A:ALA:CA	1:40:A:ALA:C	1:41:A:ASP:N	4	1.62	0.77	1.25
(1,13)	1:6:A:LEU:C	1:7:A:ASP:N	1:7:A:ASP:CA	1:7:A:ASP:C	3	3.41	1.02	3.18
(1,67)	1:64:A:ASP:C	1:65:A:ASN:N	1:65:A:ASN:CA	1:65:A:ASN:C	3	2.89	1.06	2.26
(1,135)	1:107:A:ASP:C	1:108:A:LEU:N	1:108:A:LEU:CA	1:108:A:LEU:C	3	2.36	0.95	1.98
(1,56)	1:54:A:ILE:N	1:54:A:ILE:CA	1:54:A:ILE:C	1:55:A:ILE:N	3	2.18	0.19	2.25
(1,64)	1:62:A:PHE:N	1:62:A:PHE:CA	1:62:A:PHE:C	1:63:A:GLY:N	3	2.17	1.53	1.12
(1,114)	1:95:A:ALA:N	1:95:A:ALA:CA	1:95:A:ALA:C	1:96:A:ASP:N	3	2.0	0.39	2.27
(1,112)	1:94:A:ALA:N	1:94:A:ALA:CA	1:94:A:ALA:C	1:95:A:ALA:N	3	1.89	0.48	1.95
(1,3)	1:1:A:ALA:C	1:2:A:LEU:N	1:2:A:LEU:CA	1:2:A:LEU:C	3	1.56	0.56	1.24
(1,101)	1:85:A:ASP:C	1:86:A:SER:N	1:86:A:SER:CA	1:86:A:SER:C	3	1.55	0.32	1.34
(1,26)	1:37:A:TRP:N	1:37:A:TRP:CA	1:37:A:TRP:C	1:38:A:VAL:N	3	1.52	0.59	1.14
(1,85)	1:74:A:ALA:C	1:75:A:LYS:N	1:75:A:LYS:CA	1:75:A:LYS:C	2	5.25	0.03	5.25
(1,138)	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	1:110:A:GLN:N	2	4.66	0.5	4.66
(1,118)	1:98:A:ALA:N	1:98:A:ALA:CA	1:98:A:ALA:C	1:99:A:PHE:N	2	2.92	1.63	2.92
(1,228)	1:166:A:LEU:N	1:166:A:LEU:CA	1:166:A:LEU:C	1:167:A:LEU:N	2	2.7	1.4	2.7

Continued on next page...

Continued from previous page...

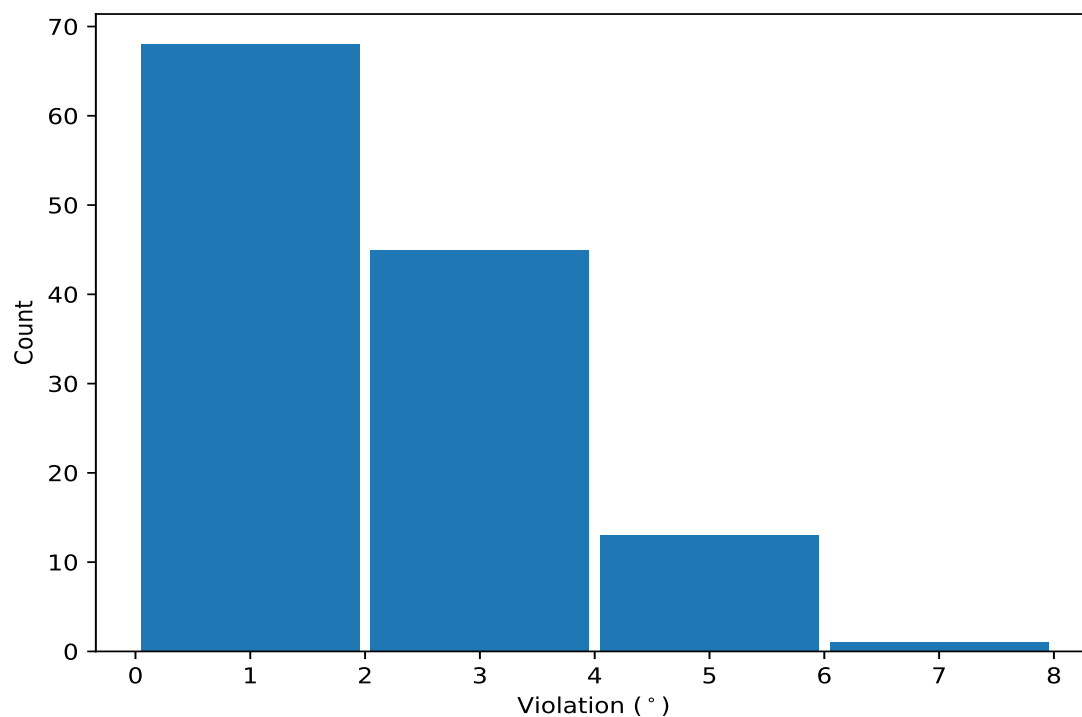
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models ¹	Mean	SD ²	Median
(1,203)	1:152:A:SER:C	1:153:A:TYR:N	1:153:A:TYR:CA	1:153:A:TYR:C	2	2.12	0.57	2.12
(1,235)	1:169:A:ALA:C	1:170:A:ALA:N	1:170:A:ALA:CA	1:170:A:ALA:C	2	2.03	0.03	2.03
(1,14)	1:7:A:ASP:N	1:7:A:ASP:CA	1:7:A:ASP:C	1:8:A:GLY:N	2	2.03	0.7	2.03
(1,54)	1:53:A:GLY:N	1:53:A:GLY:CA	1:53:A:GLY:C	1:54:A:ILE:N	2	1.94	0.43	1.94
(1,2)	1:1:A:ALA:N	1:1:A:ALA:CA	1:1:A:ALA:C	1:2:A:LEU:N	2	1.84	0.07	1.84
(1,192)	1:145:A:ASN:N	1:145:A:ASN:CA	1:145:A:ASN:C	1:146:A:THR:N	2	1.81	0.48	1.81
(1,46)	1:49:A:LEU:N	1:49:A:LEU:CA	1:49:A:LEU:C	1:50:A:ASN:N	2	1.77	0.32	1.77
(1,104)	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	1:88:A:GLY:N	2	1.71	0.16	1.71
(1,5)	1:2:A:LEU:C	1:3:A:ALA:N	1:3:A:ALA:CA	1:3:A:ALA:C	2	1.58	0.42	1.58
(1,66)	1:64:A:ASP:N	1:64:A:ASP:CA	1:64:A:ASP:C	1:65:A:ASN:N	2	1.54	0.36	1.54
(1,140)	1:110:A:GLN:N	1:110:A:GLN:CA	1:110:A:GLN:C	1:111:A:ASP:N	2	1.1	0.01	1.1

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 (°)
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	5	6.8
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	12	5.33
(1,85)	1:74:A:ALA:C	1:75:A:LYS:N	1:75:A:LYS:CA	1:75:A:LYS:C	11	5.28
(1,85)	1:74:A:ALA:C	1:75:A:LYS:N	1:75:A:LYS:CA	1:75:A:LYS:C	14	5.22
(1,138)	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	1:110:A:GLN:N	6	5.15
(1,13)	1:6:A:LEU:C	1:7:A:ASP:N	1:7:A:ASP:CA	1:7:A:ASP:C	14	4.76
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	8	4.71
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	11	4.58
(1,118)	1:98:A:ALA:N	1:98:A:ALA:CA	1:98:A:ALA:C	1:99:A:PHE:N	6	4.56
(1,67)	1:64:A:ASP:C	1:65:A:ASN:N	1:65:A:ASN:CA	1:65:A:ASN:C	6	4.38
(1,64)	1:62:A:PHE:N	1:62:A:PHE:CA	1:62:A:PHE:C	1:63:A:GLY:N	7	4.33
(1,138)	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	1:110:A:GLN:N	4	4.16
(1,228)	1:166:A:LEU:N	1:166:A:LEU:CA	1:166:A:LEU:C	1:167:A:LEU:N	1	4.1
(1,11)	1:5:A:ASP:C	1:6:A:LEU:N	1:6:A:LEU:CA	1:6:A:LEU:C	9	4.03
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	13	3.88
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	9	3.81
(1,135)	1:107:A:ASP:C	1:108:A:LEU:N	1:108:A:LEU:CA	1:108:A:LEU:C	9	3.66
(1,70)	1:66:A:THR:N	1:66:A:THR:CA	1:66:A:THR:C	1:67:A:LYS:N	7	3.56
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	14	3.38
(1,225)	1:164:A:GLY:C	1:165:A:ASP:N	1:165:A:ASP:CA	1:165:A:ASP:C	5	3.28
(1,13)	1:6:A:LEU:C	1:7:A:ASP:N	1:7:A:ASP:CA	1:7:A:ASP:C	6	3.18
(1,34)	1:41:A:ASP:N	1:41:A:ASP:CA	1:41:A:ASP:C	1:42:A:ASP:N	4	3.07
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	11	3.0
(1,68)	1:65:A:ASN:N	1:65:A:ASN:CA	1:65:A:ASN:C	1:66:A:THR:N	6	2.95
(1,32)	1:40:A:ALA:N	1:40:A:ALA:CA	1:40:A:ALA:C	1:41:A:ASP:N	9	2.94
(1,226)	1:165:A:ASP:N	1:165:A:ASP:CA	1:165:A:ASP:C	1:166:A:LEU:N	8	2.93
(1,14)	1:7:A:ASP:N	1:7:A:ASP:CA	1:7:A:ASP:C	1:8:A:GLY:N	5	2.72
(1,226)	1:165:A:ASP:N	1:165:A:ASP:CA	1:165:A:ASP:C	1:166:A:LEU:N	5	2.69
(1,203)	1:152:A:SER:C	1:153:A:TYR:N	1:153:A:TYR:CA	1:153:A:TYR:C	7	2.69
(1,62)	1:60:A:GLU:N	1:60:A:GLU:CA	1:60:A:GLU:C	1:61:A:LEU:N	3	2.55
(1,124)	1:101:A:SER:N	1:101:A:SER:CA	1:101:A:SER:C	1:102:A:LEU:N	12	2.49
(1,112)	1:94:A:ALA:N	1:94:A:ALA:CA	1:94:A:ALA:C	1:95:A:ALA:N	2	2.45
(1,134)	1:106:A:GLN:N	1:106:A:GLN:CA	1:106:A:GLN:C	1:107:A:ASP:N	12	2.41
(1,56)	1:54:A:ILE:N	1:54:A:ILE:CA	1:54:A:ILE:C	1:55:A:ILE:N	10	2.37
(1,54)	1:53:A:GLY:N	1:53:A:GLY:CA	1:53:A:GLY:C	1:54:A:ILE:N	5	2.37
(1,26)	1:37:A:TRP:N	1:37:A:TRP:CA	1:37:A:TRP:C	1:38:A:VAL:N	6	2.35
(1,3)	1:1:A:ALA:C	1:2:A:LEU:N	1:2:A:LEU:CA	1:2:A:LEU:C	2	2.34
(1,134)	1:106:A:GLN:N	1:106:A:GLN:CA	1:106:A:GLN:C	1:107:A:ASP:N	9	2.3
(1,192)	1:145:A:ASN:N	1:145:A:ASN:CA	1:145:A:ASN:C	1:146:A:THR:N	13	2.29
(1,45)	1:48:A:ASP:C	1:49:A:LEU:N	1:49:A:LEU:CA	1:49:A:LEU:C	12	2.29
(1,13)	1:6:A:LEU:C	1:7:A:ASP:N	1:7:A:ASP:CA	1:7:A:ASP:C	3	2.29
(1,114)	1:95:A:ALA:N	1:95:A:ALA:CA	1:95:A:ALA:C	1:96:A:ASP:N	9	2.28
(1,114)	1:95:A:ALA:N	1:95:A:ALA:CA	1:95:A:ALA:C	1:96:A:ASP:N	6	2.27
(1,67)	1:64:A:ASP:C	1:65:A:ASN:N	1:65:A:ASN:CA	1:65:A:ASN:C	14	2.26
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	10	2.25
(1,56)	1:54:A:ILE:N	1:54:A:ILE:CA	1:54:A:ILE:C	1:55:A:ILE:N	4	2.25
(1,133)	1:105:A:TRP:C	1:106:A:GLN:N	1:106:A:GLN:CA	1:106:A:GLN:C	7	2.2
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	3	2.15
(1,197)	1:148:A:ALA:C	1:149:A:GLN:N	1:149:A:GLN:CA	1:149:A:GLN:C	3	2.11
(1,46)	1:49:A:LEU:N	1:49:A:LEU:CA	1:49:A:LEU:C	1:50:A:ASN:N	9	2.09
(1,162)	1:123:A:GLU:N	1:123:A:GLU:CA	1:123:A:GLU:C	1:124:A:GLU:N	11	2.08

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	2	2.07
(1,235)	1:169:A:ALA:C	1:170:A:ALA:N	1:170:A:ALA:CA	1:170:A:ALA:C	14	2.06
(1,134)	1:106:A:GLN:N	1:106:A:GLN:CA	1:106:A:GLN:C	1:107:A:ASP:N	2	2.05
(1,23)	1:14:A:VAL:C	1:15:A:ALA:N	1:15:A:ALA:CA	1:15:A:ALA:C	13	2.05
(1,67)	1:64:A:ASP:C	1:65:A:ASN:N	1:65:A:ASN:CA	1:65:A:ASN:C	10	2.03
(1,45)	1:48:A:ASP:C	1:49:A:LEU:N	1:49:A:LEU:CA	1:49:A:LEU:C	10	2.02
(1,235)	1:169:A:ALA:C	1:170:A:ALA:N	1:170:A:ALA:CA	1:170:A:ALA:C	7	2.0
(1,101)	1:85:A:ASP:C	1:86:A:SER:N	1:86:A:SER:CA	1:86:A:SER:C	11	2.0
(1,5)	1:2:A:LEU:C	1:3:A:ALA:N	1:3:A:ALA:CA	1:3:A:ALA:C	5	1.99
(1,232)	1:168:A:LEU:N	1:168:A:LEU:CA	1:168:A:LEU:C	1:169:A:ALA:N	10	1.98
(1,135)	1:107:A:ASP:C	1:108:A:LEU:N	1:108:A:LEU:CA	1:108:A:LEU:C	2	1.98
(1,112)	1:94:A:ALA:N	1:94:A:ALA:CA	1:94:A:ALA:C	1:95:A:ALA:N	12	1.95
(1,11)	1:5:A:ASP:C	1:6:A:LEU:N	1:6:A:LEU:CA	1:6:A:LEU:C	10	1.95
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	14	1.92
(1,120)	1:99:A:PHE:N	1:99:A:PHE:CA	1:99:A:PHE:C	1:100:A:GLN:N	5	1.91
(1,56)	1:54:A:ILE:N	1:54:A:ILE:CA	1:54:A:ILE:C	1:55:A:ILE:N	14	1.91
(1,2)	1:1:A:ALA:N	1:1:A:ALA:CA	1:1:A:ALA:C	1:2:A:LEU:N	12	1.91
(1,134)	1:106:A:GLN:N	1:106:A:GLN:CA	1:106:A:GLN:C	1:107:A:ASP:N	5	1.9
(1,66)	1:64:A:ASP:N	1:64:A:ASP:CA	1:64:A:ASP:C	1:65:A:ASN:N	5	1.9
(1,104)	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	1:88:A:GLY:N	14	1.87
(1,2)	1:1:A:ALA:N	1:1:A:ALA:CA	1:1:A:ALA:C	1:2:A:LEU:N	11	1.76
(1,229)	1:166:A:LEU:C	1:167:A:LEU:N	1:167:A:LEU:CA	1:167:A:LEU:C	3	1.75
(1,11)	1:5:A:ASP:C	1:6:A:LEU:N	1:6:A:LEU:CA	1:6:A:LEU:C	12	1.72
(1,11)	1:5:A:ASP:C	1:6:A:LEU:N	1:6:A:LEU:CA	1:6:A:LEU:C	7	1.68
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	1	1.62
(1,18)	1:11:A:ILE:N	1:11:A:ILE:CA	1:11:A:ILE:C	1:12:A:GLU:N	14	1.57
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	7	1.56
(1,104)	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	1:88:A:GLY:N	13	1.55
(1,203)	1:152:A:SER:C	1:153:A:TYR:N	1:153:A:TYR:CA	1:153:A:TYR:C	9	1.54
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	8	1.53
(1,54)	1:53:A:GLY:N	1:53:A:GLY:CA	1:53:A:GLY:C	1:54:A:ILE:N	8	1.52
(1,175)	1:130:A:LEU:C	1:131:A:ASP:N	1:131:A:ASP:CA	1:131:A:ASP:C	13	1.49
(1,114)	1:95:A:ALA:N	1:95:A:ALA:CA	1:95:A:ALA:C	1:96:A:ASP:N	3	1.45
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	6	1.45
(1,46)	1:49:A:LEU:N	1:49:A:LEU:CA	1:49:A:LEU:C	1:50:A:ASN:N	3	1.45
(1,135)	1:107:A:ASP:C	1:108:A:LEU:N	1:108:A:LEU:CA	1:108:A:LEU:C	1	1.44
(1,226)	1:165:A:ASP:N	1:165:A:ASP:CA	1:165:A:ASP:C	1:166:A:LEU:N	7	1.43
(1,103)	1:86:A:SER:C	1:87:A:ASN:N	1:87:A:ASN:CA	1:87:A:ASN:C	13	1.38
(1,137)	1:108:A:LEU:C	1:109:A:ASN:N	1:109:A:ASN:CA	1:109:A:ASN:C	6	1.36
(1,32)	1:40:A:ALA:N	1:40:A:ALA:CA	1:40:A:ALA:C	1:41:A:ASP:N	12	1.36
(1,101)	1:85:A:ASP:C	1:86:A:SER:N	1:86:A:SER:CA	1:86:A:SER:C	8	1.34
(1,14)	1:7:A:ASP:N	1:7:A:ASP:CA	1:7:A:ASP:C	1:8:A:GLY:N	2	1.33
(1,192)	1:145:A:ASN:N	1:145:A:ASN:CA	1:145:A:ASN:C	1:146:A:THR:N	5	1.32
(1,101)	1:85:A:ASP:C	1:86:A:SER:N	1:86:A:SER:CA	1:86:A:SER:C	3	1.32
(1,134)	1:106:A:GLN:N	1:106:A:GLN:CA	1:106:A:GLN:C	1:107:A:ASP:N	11	1.31
(1,134)	1:106:A:GLN:N	1:106:A:GLN:CA	1:106:A:GLN:C	1:107:A:ASP:N	8	1.3
(1,228)	1:166:A:LEU:N	1:166:A:LEU:CA	1:166:A:LEU:C	1:167:A:LEU:N	6	1.29
(1,226)	1:165:A:ASP:N	1:165:A:ASP:CA	1:165:A:ASP:C	1:166:A:LEU:N	6	1.29
(1,118)	1:98:A:ALA:N	1:98:A:ALA:CA	1:98:A:ALA:C	1:99:A:PHE:N	9	1.29
(1,112)	1:94:A:ALA:N	1:94:A:ALA:CA	1:94:A:ALA:C	1:95:A:ALA:N	5	1.27
(1,11)	1:5:A:ASP:C	1:6:A:LEU:N	1:6:A:LEU:CA	1:6:A:LEU:C	13	1.27

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,195)	1:146:A:THR:C	1:147:A:LEU:N	1:147:A:LEU:CA	1:147:A:LEU:C	10	1.26
(1,3)	1:1:A:ALA:C	1:2:A:LEU:N	1:2:A:LEU:CA	1:2:A:LEU:C	1	1.24
(1,60)	1:59:A:ALA:N	1:59:A:ALA:CA	1:59:A:ALA:C	1:60:A:GLU:N	10	1.22
(1,79)	1:70:A:ASP:C	1:71:A:GLY:N	1:71:A:GLY:CA	1:71:A:GLY:C	9	1.21
(1,15)	1:7:A:ASP:C	1:8:A:GLY:N	1:8:A:GLY:CA	1:8:A:GLY:C	13	1.21
(1,207)	1:154:A:THR:C	1:155:A:LYS:N	1:155:A:LYS:CA	1:155:A:LYS:C	2	1.19
(1,66)	1:64:A:ASP:N	1:64:A:ASP:CA	1:64:A:ASP:C	1:65:A:ASN:N	10	1.19
(1,45)	1:48:A:ASP:C	1:49:A:LEU:N	1:49:A:LEU:CA	1:49:A:LEU:C	2	1.17
(1,5)	1:2:A:LEU:C	1:3:A:ALA:N	1:3:A:ALA:CA	1:3:A:ALA:C	13	1.16
(1,123)	1:100:A:GLN:C	1:101:A:SER:N	1:101:A:SER:CA	1:101:A:SER:C	13	1.15
(1,58)	1:55:A:ILE:N	1:55:A:ILE:CA	1:55:A:ILE:C	1:56:A:ASP:N	13	1.15
(1,32)	1:40:A:ALA:N	1:40:A:ALA:CA	1:40:A:ALA:C	1:41:A:ASP:N	14	1.14
(1,26)	1:37:A:TRP:N	1:37:A:TRP:CA	1:37:A:TRP:C	1:38:A:VAL:N	13	1.14
(1,64)	1:62:A:PHE:N	1:62:A:PHE:CA	1:62:A:PHE:C	1:63:A:GLY:N	5	1.12
(1,185)	1:139:A:LYS:C	1:140:A:ASN:N	1:140:A:ASN:CA	1:140:A:ASN:C	12	1.11
(1,140)	1:110:A:GLN:N	1:110:A:GLN:CA	1:110:A:GLN:C	1:111:A:ASP:N	8	1.1
(1,140)	1:110:A:GLN:N	1:110:A:GLN:CA	1:110:A:GLN:C	1:111:A:ASP:N	4	1.09
(1,3)	1:1:A:ALA:C	1:2:A:LEU:N	1:2:A:LEU:CA	1:2:A:LEU:C	5	1.09
(1,64)	1:62:A:PHE:N	1:62:A:PHE:CA	1:62:A:PHE:C	1:63:A:GLY:N	10	1.07
(1,50)	1:51:A:GLY:N	1:51:A:GLY:CA	1:51:A:GLY:C	1:52:A:ASN:N	4	1.07
(1,45)	1:48:A:ASP:C	1:49:A:LEU:N	1:49:A:LEU:CA	1:49:A:LEU:C	9	1.06
(1,26)	1:37:A:TRP:N	1:37:A:TRP:CA	1:37:A:TRP:C	1:38:A:VAL:N	7	1.06
(1,4)	1:2:A:LEU:N	1:2:A:LEU:CA	1:2:A:LEU:C	1:3:A:ALA:N	2	1.04
(1,33)	1:40:A:ALA:C	1:41:A:ASP:N	1:41:A:ASP:CA	1:41:A:ASP:C	12	1.02
(1,32)	1:40:A:ALA:N	1:40:A:ALA:CA	1:40:A:ALA:C	1:41:A:ASP:N	5	1.02