Full wwPDB NMR Structure Validation Report (i)

Nov 28, 2023 – 04:21 PM EST

PDB ID	:	2LFK
BMRB ID	:	17763
Title	:	NMR solution structure of native TdPI-short
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Deposited on	:	2011-07-06

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

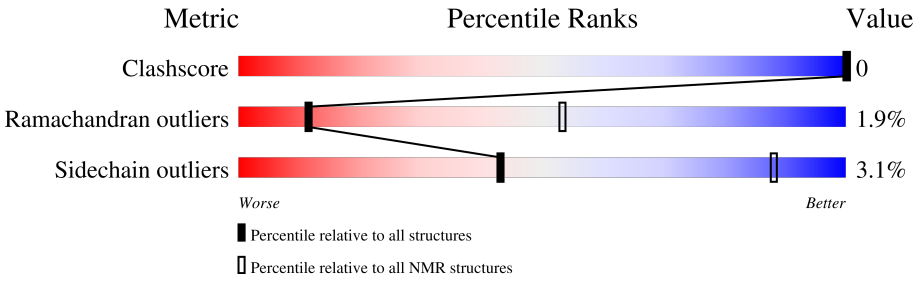
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
BMRB Restraints Analysis	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
SOLUTION NMR

The overall completeness of chemical shifts assignment is 49%.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for >=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 <=5%

Mol	Chain	Length	Quality of chain
1	A	57	82% • 16%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 8 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *fewest violations*.

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

Well-defined (core) protein residues			
Well-defined core	**Residue range (total)**	**Backbone RMSD (Å)**	**Medoid model**
1	A:26-A:73 (48)	0.69	8

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. No single-model clusters were found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 840 atoms, of which 404 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Tryptase inhibitor.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	57	840	268	404	76	82	10	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	GLY	-	expression tag	UNP Q1EG59
A	20	ASP	-	expression tag	UNP Q1EG59

4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

• Molecule 1: Tryptase inhibitor

Chain A: 82% • 16%

G19 D20 K21 E22 E23 C24 T25 G57 P74 G75

4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: Tryptase inhibitor

Chain A: 75% 9% 16%

G19 D20 K21 E22 E23 C24 T25 E32 K39 G57 T60 C73 P74 G75

4.2.2 Score per residue for model 2

• Molecule 1: Tryptase inhibitor

Chain A: 82% • 16%

G19 D20 K21 E22 E23 C24 T25 C73 P74 G75

4.2.3 Score per residue for model 3

- Molecule 1: Tryptase inhibitor

Chain A: 81% • 16%

G19 D20 K21 E22 E23 C24 T25 G57 T60 P74 G75

4.2.4 Score per residue for model 4

- Molecule 1: Tryptase inhibitor

Chain A: 84% 16%

G19 D20 K21 E22 E23 C24 T25 P74 G75

4.2.5 Score per residue for model 5

- Molecule 1: Tryptase inhibitor

Chain A: 79% 5% 16%

G19 D20 K21 E22 E23 C24 T25 E32 K39 C51 P74 G75

4.2.6 Score per residue for model 6

- Molecule 1: Tryptase inhibitor

Chain A: 81% • 16%

G19 D20 K21 E22 E23 C24 T25 K39 G57 P74 G75

4.2.7 Score per residue for model 7

- Molecule 1: Tryptase inhibitor

Chain A: 81% • 16%

G19 D20 K21 E22 E23 C24 T25 G57 R71 P74 G75

4.2.8 Score per residue for model 8 (medoid)

- Molecule 1: Tryptase inhibitor

Chain A: 82% • 16%

G19 D20 K21 E22 E23 C24 T25 G57 P74 G75

4.2.9 Score per residue for model 9

- Molecule 1: Tryptase inhibitor

Chain A: 84% 16%

G19 D20 K21 E22 E23 C24 T25 P74 G75

4.2.10 Score per residue for model 10

- Molecule 1: Tryptase inhibitor

Chain A: 77% 7% 16%

G19 D20 K21 E22 E23 C24 T25 E32 E56 G57 C73 P74 G75

4.2.11 Score per residue for model 11

- Molecule 1: Tryptase inhibitor

Chain A: 81% • 16%

G19 D20 K21 E22 E23 C24 T25 T60 R65 P74 G75

4.2.12 Score per residue for model 12

- Molecule 1: Tryptase inhibitor

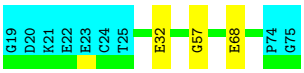
Chain A: 81% • 16%

G19 D20 K21 E22 E23 C24 T25 G57 C73 P74 G75

4.2.13 Score per residue for model 13

- Molecule 1: Tryptase inhibitor

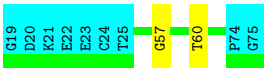
Chain A: 79% 5% 16%



4.2.14 Score per residue for model 14

- Molecule 1: Tryptase inhibitor

Chain A: 81% • 16%



4.2.15 Score per residue for model 15

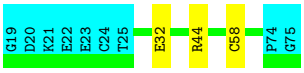
- Molecule 1: Tryptase inhibitor

Chain A: 77% 7% 16%

4.2.16 Score per residue for model 16

- Molecule 1: Tryptase inhibitor

Chain A: 79% 5% 16%



4.2.17 Score per residue for model 17

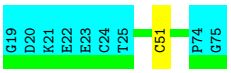
- Molecule 1: Tryptase inhibitor

Chain A: 75% 9% 16%

4.2.18 Score per residue for model 18

- Molecule 1: Tryptase inhibitor

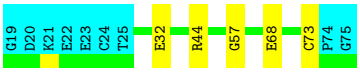
Chain A:



4.2.19 Score per residue for model 19

- Molecule 1: Tryptase inhibitor

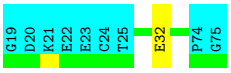
Chain A:



4.2.20 Score per residue for model 20

- Molecule 1: Tryptase inhibitor

Chain A:



5 Refinement protocol and experimental data overview (i)

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

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

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

Software name	Classification	Version
CYANA	structure solution	2.1
Amber	refinement	9.0
CYANA	refinement	

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

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	356
Number of shifts mapped to atoms	356
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	49%

6 Model quality (i)

6.1 Standard geometry (i)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	0.84±0.01	0±0/382 (0.0± 0.0%)	1.00±0.02	0±1/514 (0.1± 0.1%)
All	All	0.84	0/7640 (0.0%)	1.00	6/10280 (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(^{o})	Ideal(^{o})	Models	
								Worst	Total
1	A	71	ARG	NE-CZ-NH1	6.31	123.45	120.30	7	3
1	A	41	ARG	NE-CZ-NH1	5.65	123.12	120.30	15	1
1	A	44	ARG	NE-CZ-NH1	5.26	122.93	120.30	19	2

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
All	All	7440	6980	6980	-

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 -.

There are no clashes.

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	48/57 (84%)	44±2 (92±4%)	3±2 (6±3%)	1±1 (2±1%)	11 53
All	All	960/1140 (84%)	888 (92%)	54 (6%)	18 (2%)	11 53

All 3 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	57	GLY	12
1	A	60	THR	5
1	A	58	CYS	1

6.3.2 Protein sidechains (i)

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	39/46 (85%)	38±1 (97±3%)	1±1 (3±3%)	43 88
All	All	780/920 (85%)	756 (97%)	24 (3%)	43 88

All 8 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	32	GLU	9
1	A	73	CYS	5
1	A	39	LYS	3
1	A	51	CYS	2
1	A	68	GLU	2
1	A	56	GLU	1
1	A	65	ARG	1

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Mol	Chain	Res	Type	Models (Total)
1	A	48	MET	1

6.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

6.5 Carbohydrates (i)

There are no monosaccharides in this entry.

6.6 Ligand geometry (i)

There are no ligands in this entry.

6.7 Other polymers (i)

There are no such molecules in this entry.

6.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

7 Chemical shift validation (i)

The completeness of assignment taking into account all chemical shift lists is 49% for the well-defined parts and 49% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping (i)

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

Total number of shifts	356
Number of shifts mapped to atoms	356
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	2

7.1.2 Chemical shift referencing (i)

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

7.1.3 Completeness of resonance assignments (i)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 49%, i.e. 305 atoms were assigned a chemical shift out of a possible 618. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	**^{1}H**	**^{13}C**	**^{15}N**
Backbone	100/243 (41%)	100/101 (99%)	0/96 (0%)	0/46 (0%)
Sidechain	176/308 (57%)	176/199 (88%)	0/92 (0%)	0/17 (0%)
Aromatic	29/67 (43%)	29/31 (94%)	0/35 (0%)	0/1 (0%)
Overall	305/618 (49%)	305/331 (92%)	0/223 (0%)	0/64 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 49%, i.e. 350 atoms were assigned a chemical shift out of a possible 712. 0 out of 4 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	117/288 (41%)	117/120 (98%)	0/114 (0%)	0/54 (0%)
Sidechain	204/357 (57%)	204/229 (89%)	0/110 (0%)	0/18 (0%)
Aromatic	29/67 (43%)	29/31 (94%)	0/35 (0%)	0/1 (0%)
Overall	350/712 (49%)	350/380 (92%)	0/259 (0%)	0/73 (0%)

7.1.4 Statistically unusual chemical shifts (i)

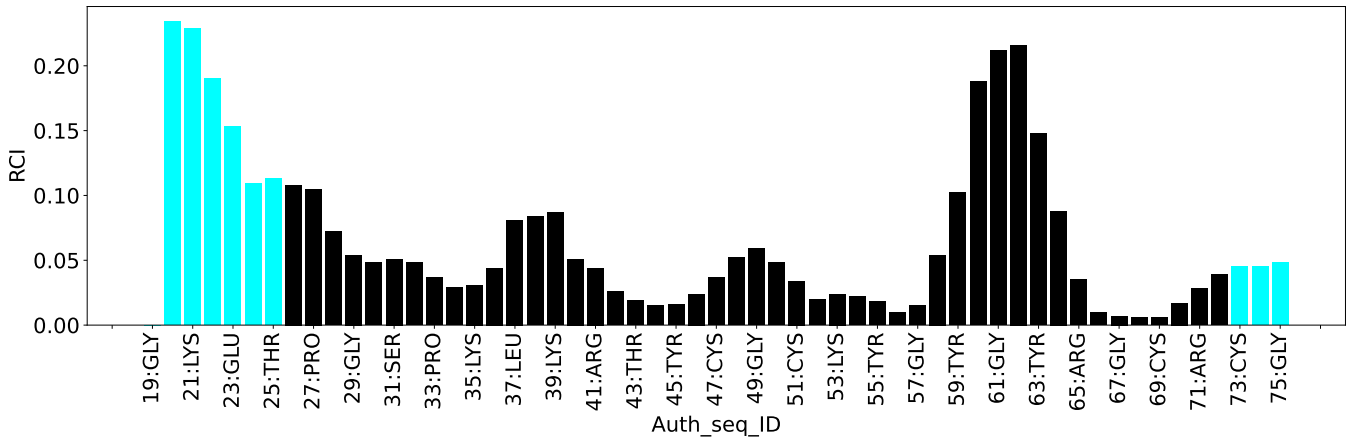
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
1	A	33	PRO	HG3	0.07	0.33 – 3.48	-5.8
1	A	33	PRO	HG2	0.20	0.41 – 3.45	-5.7

7.1.5 Random Coil Index (RCI) plots (i)

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

Random coil index (RCI) for chain A:



8 NMR restraints analysis (i)

8.1 Conformationally restricting restraints (i)

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

Description	Value
Total distance restraints	693
Intra-residue (|i-j|=0)	200
Sequential (|i-j|=1)	181
Medium range (|i-j|>1 and |i-j|<5)	115
Long range (|i-j|≥5)	193
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	4
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	12.2
Number of long range restraints per residue[1]	3.5

[1]Long range hydrogen bonds and disulfide bonds are counted as long range restraints while calculating the number of long range restraints per residue

8.2 Residual restraint violations (i)

This section provides the overview of the restraint violations analysis. The violations are binned as small, medium and large violations based on its absolute value. Average number of violations per model is calculated by dividing the total number of violations in each bin by the size of the ensemble.

8.2.1 Average number of distance violations per model (i)

Distance violations less than 0.1 Å are not included in the calculation.

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	1.6	0.2
0.2-0.5 (Medium)	0.7	0.41
>0.5 (Large)	2.0	1.34

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

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

9 Distance violation analysis (i)

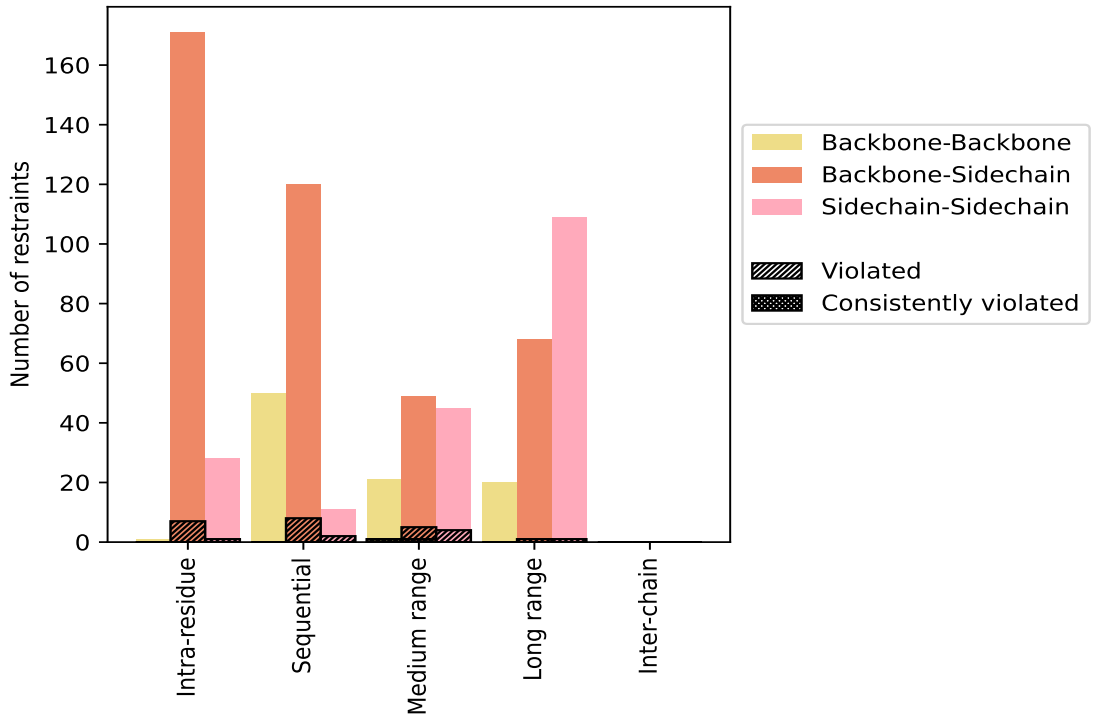
9.1 Summary of distance violations (i)

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

Restraints type	Count	%[1]	Violated[3]			Consistently Violated[4]		
			Count	%[2]	%[1]	Count	%[2]	%[1]
Intra-residue (|i-j|=0)	200	28.9	8	4.0	1.2	0	0.0	0.0
Backbone-Backbone	1	0.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	171	24.7	7	4.1	1.0	0	0.0	0.0
Sidechain-Sidechain	28	4.0	1	3.6	0.1	0	0.0	0.0
Sequential (|i-j|=1)	181	26.1	10	5.5	1.4	0	0.0	0.0
Backbone-Backbone	50	7.2	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	120	17.3	8	6.7	1.2	0	0.0	0.0
Sidechain-Sidechain	11	1.6	2	18.2	0.3	0	0.0	0.0
Medium range (|i-j|>1 & |i-j|<5)	115	16.6	10	8.7	1.4	2	1.7	0.3
Backbone-Backbone	21	3.0	1	4.8	0.1	1	4.8	0.1
Backbone-Sidechain	49	7.1	5	10.2	0.7	1	2.0	0.1
Sidechain-Sidechain	45	6.5	4	8.9	0.6	0	0.0	0.0
Long range (|i-j|≥5)	193	27.8	2	1.0	0.3	0	0.0	0.0
Backbone-Backbone	20	2.9	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	68	9.8	1	1.5	0.1	0	0.0	0.0
Sidechain-Sidechain	105	15.2	1	1.0	0.1	0	0.0	0.0
Inter-chain	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Hydrogen bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	4	0.6	0	0.0	0.0	0	0.0	0.0
Total	693	100.0	30	4.3	4.3	2	0.3	0.3
Backbone-Backbone	92	13.3	1	1.1	0.1	1	1.1	0.1
Backbone-Sidechain	408	58.9	21	5.1	3.0	1	0.2	0.1
Sidechain-Sidechain	193	27.8	8	4.1	1.2	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	1	2	0	0	3	0.76	1.13	0.46	1.04
2	1	0	2	0	0	3	0.6	0.9	0.33	0.77
3	0	0	2	0	0	2	0.84	0.89	0.05	0.84
4	1	0	3	0	0	4	0.48	0.89	0.36	0.48
5	1	2	5	0	0	8	0.34	0.85	0.28	0.2
6	1	1	4	0	0	6	0.5	0.86	0.3	0.57
7	1	1	2	0	0	4	0.65	1.17	0.48	0.65
8	0	0	2	0	0	2	0.92	0.94	0.02	0.92
9	0	0	2	0	0	2	1.04	1.05	0.02	1.04
10	0	1	2	0	0	3	0.6	0.88	0.33	0.79
11	1	2	2	0	0	5	0.4	0.82	0.31	0.19

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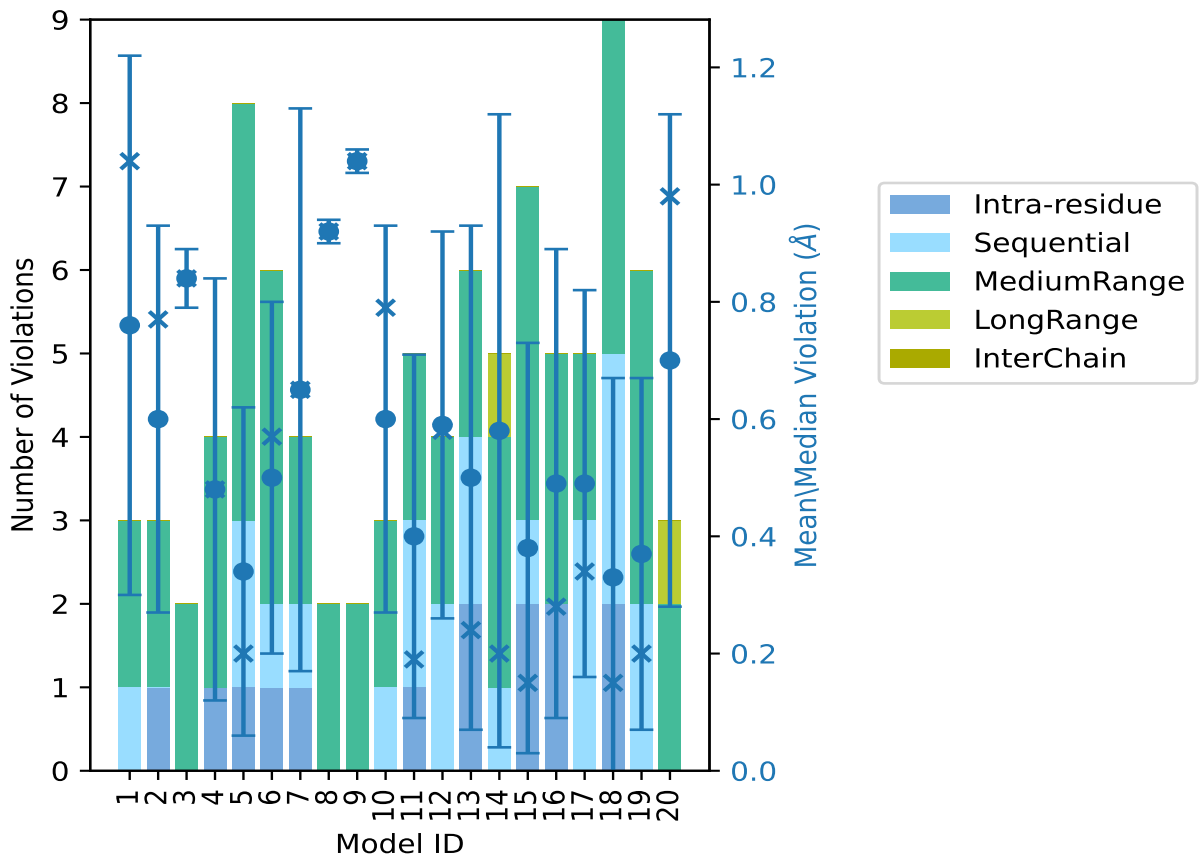
Model ID	Number of violations						Mean (Å)	Max (Å)	SD[6] (Å)	Median (Å)
	IR[1]	SQ[2]	MR[3]	LR[4]	IC[5]	Total				
12	0	2	2	0	0	4	0.59	0.95	0.33	0.58
13	2	2	2	0	0	6	0.5	1.12	0.43	0.24
14	0	1	3	1	0	5	0.58	1.34	0.54	0.2
15	2	1	4	0	0	7	0.38	1.01	0.35	0.15
16	2	0	3	0	0	5	0.49	1.03	0.4	0.28
17	0	3	2	0	0	5	0.49	0.91	0.33	0.34
18	2	3	4	0	0	9	0.33	1.04	0.34	0.15
19	0	2	4	0	0	6	0.37	0.8	0.3	0.2
20	0	0	2	1	0	3	0.7	1.02	0.42	0.98

[1]Intra-residue restraints, [2]Sequential restraints, [3]Medium range restraints, [4]Long range restraints, [5]Inter-chain restraints, [6]Standard deviation

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



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

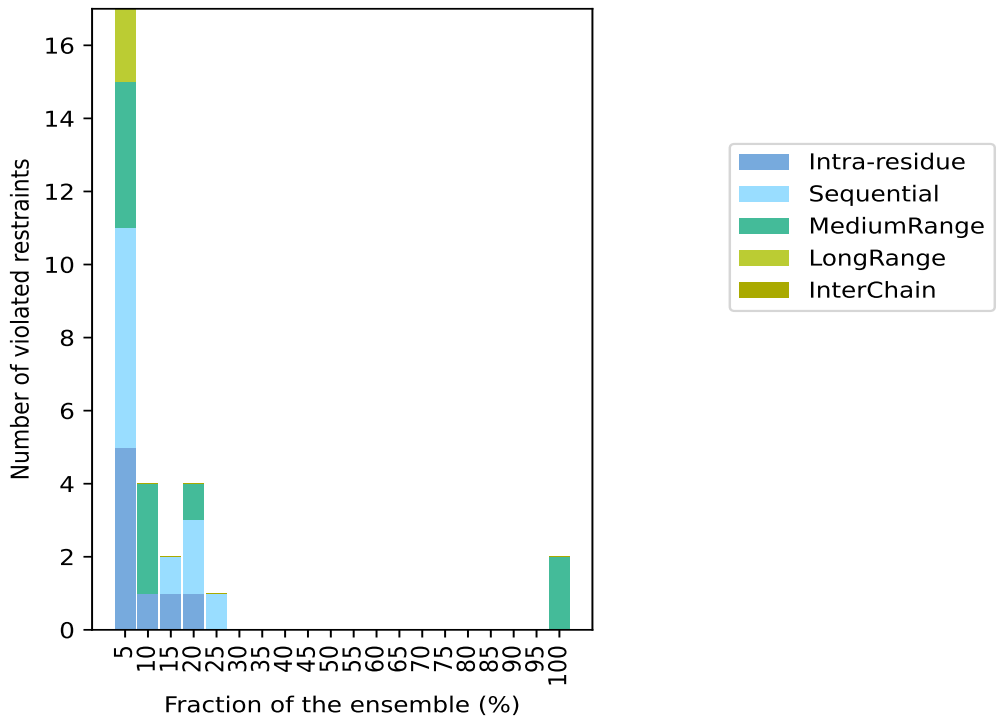
9.3 Distance violation statistics for the ensemble (i)

Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 659(IR:192, SQ:171, MR:105, LR:191, 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]	%
5	6	4	2	0	17	1	5.0
1	0	3	0	0	4	2	10.0
1	1	0	0	0	2	3	15.0
1	2	1	0	0	4	4	20.0
0	1	0	0	0	1	5	25.0
0	0	0	0	0	0	6	30.0
0	0	0	0	0	0	7	35.0
0	0	0	0	0	0	8	40.0
0	0	0	0	0	0	9	45.0
0	0	0	0	0	0	10	50.0
0	0	0	0	0	0	11	55.0
0	0	0	0	0	0	12	60.0
0	0	0	0	0	0	13	65.0
0	0	0	0	0	0	14	70.0
0	0	0	0	0	0	15	75.0
0	0	0	0	0	0	16	80.0
0	0	0	0	0	0	17	85.0
0	0	0	0	0	0	18	90.0
0	0	0	0	0	0	19	95.0
0	0	2	0	0	2	20	100.0

[1]Intra-residue restraints, [2]Sequential restraints, [3]Medium range restraints, [4]Long range restraints, [5]Inter-chain restraints, [6] Number of models with violations

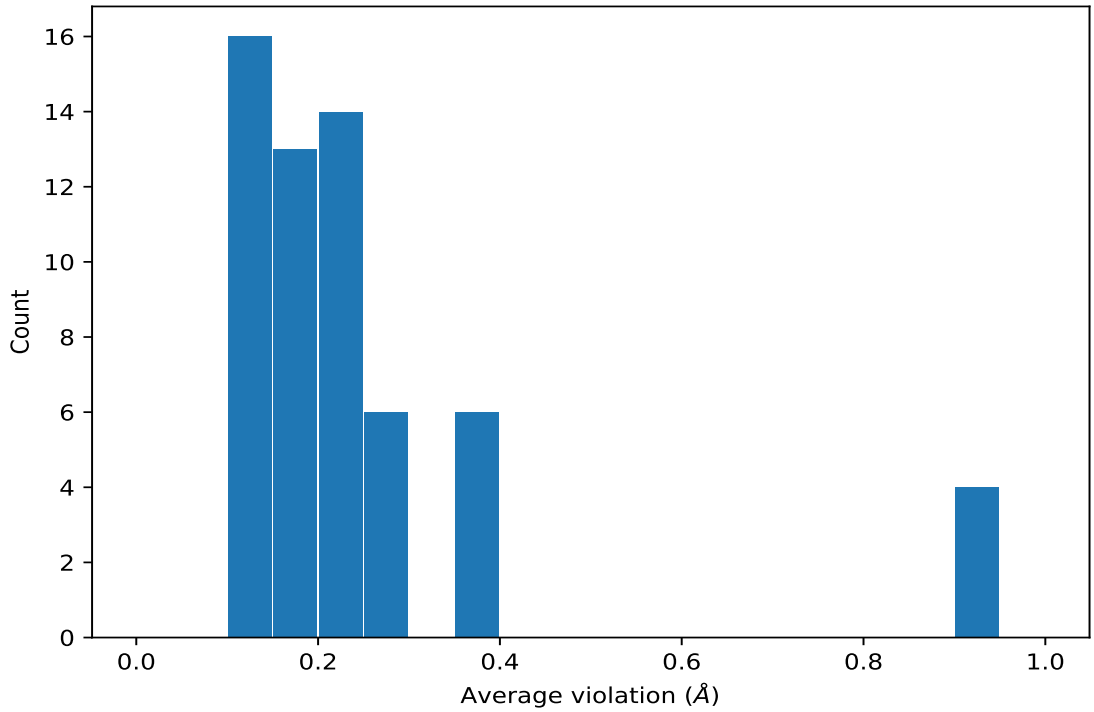
9.3.1 Bar graph : Distance violation statistics for the ensemble (i)



9.4 Most violated distance restraints in the ensemble (i)

9.4.1 Histogram : Distribution of mean distance violations (i)

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



9.4.2 Table: Most violated distance restraints (i)

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

Key	Atom-1	Atom-2	Models[1]	Mean (Å)	SD[1] (Å)	Median (Å)
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	20	0.95	0.14	0.9
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	20	0.95	0.14	0.9
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	20	0.95	0.14	0.9
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	20	0.92	0.13	0.89
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD11	5	0.39	0.19	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD12	5	0.39	0.19	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD13	5	0.39	0.19	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD21	5	0.39	0.19	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD22	5	0.39	0.19	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD23	5	0.39	0.19	0.33
(1,661)	1:66:A:MET:HG2	1:67:A:GLY:HA3	4	0.21	0.05	0.23
(1,661)	1:66:A:MET:HG3	1:67:A:GLY:HA3	4	0.21	0.05	0.23
(1,175)	1:26:A:VAL:HG11	1:30:A:TRP:HD1	4	0.19	0.04	0.18
(1,175)	1:26:A:VAL:HG12	1:30:A:TRP:HD1	4	0.19	0.04	0.18
(1,175)	1:26:A:VAL:HG13	1:30:A:TRP:HD1	4	0.19	0.04	0.18
(1,175)	1:26:A:VAL:HG21	1:30:A:TRP:HD1	4	0.19	0.04	0.18

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Key	Atom-1	Atom-2	Models[1]	Mean (Å)	SD[1] (Å)	Median (Å)
(1,175)	1:26:A:VAL:HG22	1:30:A:TRP:HD1	4	0.19	0.04	0.18
(1,175)	1:26:A:VAL:HG23	1:30:A:TRP:HD1	4	0.19	0.04	0.18
(1,340)	1:26:A:VAL:HG11	1:27:A:PRO:HD3	4	0.16	0.05	0.15
(1,340)	1:26:A:VAL:HG12	1:27:A:PRO:HD3	4	0.16	0.05	0.15
(1,340)	1:26:A:VAL:HG13	1:27:A:PRO:HD3	4	0.16	0.05	0.15
(1,340)	1:26:A:VAL:HG21	1:27:A:PRO:HD3	4	0.16	0.05	0.15
(1,340)	1:26:A:VAL:HG22	1:27:A:PRO:HD3	4	0.16	0.05	0.15
(1,340)	1:26:A:VAL:HG23	1:27:A:PRO:HD3	4	0.16	0.05	0.15
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG2	4	0.14	0.02	0.15
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG3	4	0.14	0.02	0.15
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG11	3	0.27	0.05	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG12	3	0.27	0.05	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG13	3	0.27	0.05	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG21	3	0.27	0.05	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG22	3	0.27	0.05	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG23	3	0.27	0.05	0.27
(1,178)	1:44:A:ARG:HA	1:44:A:ARG:HE	3	0.17	0.02	0.19
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD11	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD12	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD13	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD21	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD22	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD23	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD11	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD12	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD13	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD21	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD22	2	0.2	0.08	0.2
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD23	2	0.2	0.08	0.2
(1,87)	1:23:A:GLU:H	1:23:A:GLU:HG2	2	0.15	0.0	0.15
(1,87)	1:23:A:GLU:H	1:23:A:GLU:HG3	2	0.15	0.0	0.15
(1,454)	1:26:A:VAL:HG11	1:30:A:TRP:HE3	2	0.12	0.01	0.12
(1,454)	1:26:A:VAL:HG12	1:30:A:TRP:HE3	2	0.12	0.01	0.12
(1,454)	1:26:A:VAL:HG13	1:30:A:TRP:HE3	2	0.12	0.01	0.12
(1,454)	1:26:A:VAL:HG21	1:30:A:TRP:HE3	2	0.12	0.01	0.12
(1,454)	1:26:A:VAL:HG22	1:30:A:TRP:HE3	2	0.12	0.01	0.12
(1,454)	1:26:A:VAL:HG23	1:30:A:TRP:HE3	2	0.12	0.01	0.12
(1,458)	1:26:A:VAL:HG11	1:30:A:TRP:HB2	2	0.12	0.01	0.12
(1,458)	1:26:A:VAL:HG12	1:30:A:TRP:HB2	2	0.12	0.01	0.12
(1,458)	1:26:A:VAL:HG13	1:30:A:TRP:HB2	2	0.12	0.01	0.12
(1,458)	1:26:A:VAL:HG21	1:30:A:TRP:HB2	2	0.12	0.01	0.12
(1,458)	1:26:A:VAL:HG22	1:30:A:TRP:HB2	2	0.12	0.01	0.12

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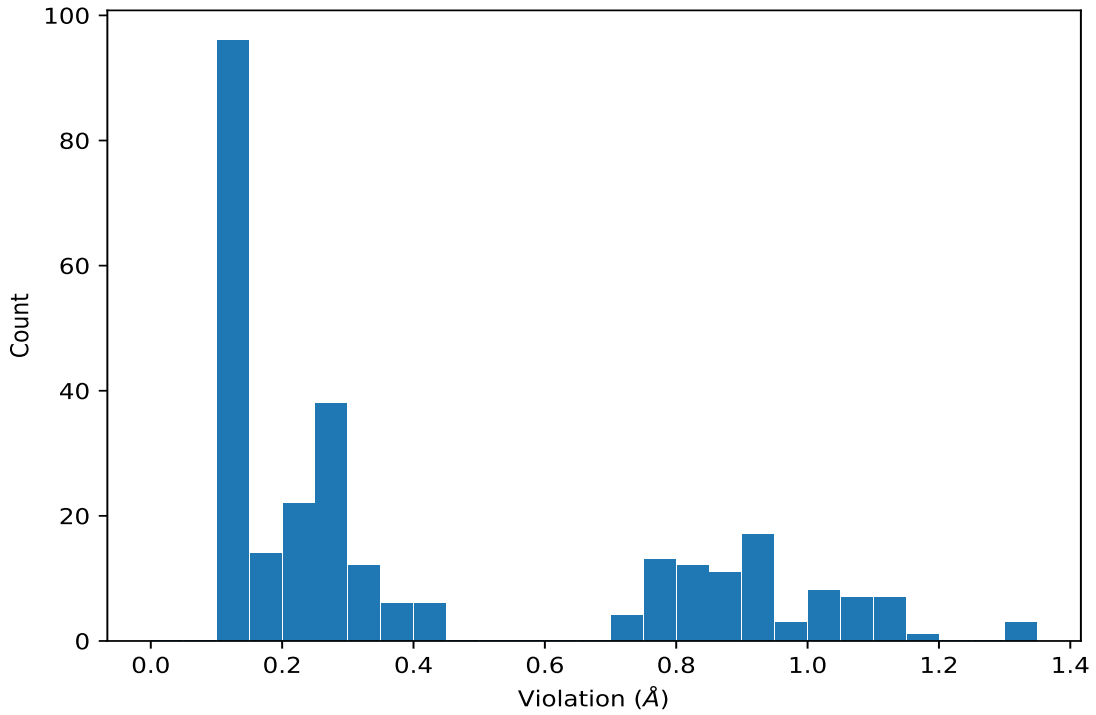
Key	Atom-1	Atom-2	Models[1]	Mean (Å)	SD[1] (Å)	Median (Å)
(1,458)	1:26:A:VAL:HG23	1:30:A:TRP:HB2	2	0.12	0.01	0.12

[1]Number of violated models, [2]Standard deviation

9.5 All violated distance restraints (i)

9.5.1 Histogram : Distribution of distance violations (i)

The following histogram shows the distribution of the absolute value of the violation for all violated restraints in the ensemble.



9.5.2 Table : All distance violations (i)

The following table lists the absolute value of the violation for each restraint in the ensemble sorted by its value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	14	1.34
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	14	1.34
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	14	1.34

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	7	1.17
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	1	1.13
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	1	1.13
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	1	1.13
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	13	1.12
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	13	1.12
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	13	1.12
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	14	1.11
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	7	1.09
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	7	1.09
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	7	1.09
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	13	1.09
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	9	1.05
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	9	1.05
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	9	1.05
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	1	1.04
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	18	1.04
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	16	1.03
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	9	1.02
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	20	1.02
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	15	1.01
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	15	1.01
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	15	1.01
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	20	0.98
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	20	0.98
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	20	0.98
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	12	0.95
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	8	0.94
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	8	0.94
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	8	0.94
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	16	0.91
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	16	0.91
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	16	0.91
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	17	0.91
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	17	0.91
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	17	0.91
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	2	0.9
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	2	0.9
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	2	0.9
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	12	0.9
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	12	0.9
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	12	0.9

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	8	0.9
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	3	0.89
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	3	0.89
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	3	0.89
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	4	0.89
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	4	0.89
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	4	0.89
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	18	0.89
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	18	0.89
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	18	0.89
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	10	0.88
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	6	0.86
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	5	0.85
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	5	0.85
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	5	0.85
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	17	0.85
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	11	0.82
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	11	0.82
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	11	0.82
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	19	0.8
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	19	0.8
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	19	0.8
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	4	0.8
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	15	0.8
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	10	0.79
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	10	0.79
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	10	0.79
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	3	0.78
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	5	0.78
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	19	0.78
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	2	0.77
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD11	6	0.76
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD12	6	0.76
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD13	6	0.76
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD21	6	0.76
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD22	6	0.76
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD23	6	0.76
(1,59)	1:67:A:GLY:HA3	1:70:A:ALA:H	11	0.74
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB1	6	0.73
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB2	6	0.73
(1,436)	1:67:A:GLY:HA3	1:70:A:ALA:HB3	6	0.73
(1,554)	1:35:A:LYS:HA	1:37:A:LEU:HD11	6	0.41

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,554)	1:35:A:LYS:HA	1:37:A:LEU:HD12	6	0.41
(1,554)	1:35:A:LYS:HA	1:37:A:LEU:HD13	6	0.41
(1,554)	1:35:A:LYS:HA	1:37:A:LEU:HD21	6	0.41
(1,554)	1:35:A:LYS:HA	1:37:A:LEU:HD22	6	0.41
(1,554)	1:35:A:LYS:HA	1:37:A:LEU:HD23	6	0.41
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD11	5	0.36
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD12	5	0.36
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD13	5	0.36
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD21	5	0.36
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD22	5	0.36
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD23	5	0.36
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG11	17	0.34
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG12	17	0.34
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG13	17	0.34
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG21	17	0.34
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG22	17	0.34
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG23	17	0.34
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD11	15	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD12	15	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD13	15	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD21	15	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD22	15	0.33
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD23	15	0.33
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD11	16	0.28
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD12	16	0.28
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD13	16	0.28
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD21	16	0.28
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD22	16	0.28
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD23	16	0.28
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD11	16	0.28
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD12	16	0.28
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD13	16	0.28
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD21	16	0.28
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD22	16	0.28
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD23	16	0.28
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG11	12	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG12	12	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG13	12	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG21	12	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG22	12	0.27
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG23	12	0.27
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD11	19	0.26

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD12	19	0.26
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD13	19	0.26
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD21	19	0.26
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD22	19	0.26
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD23	19	0.26
(1,661)	1:66:A:MET:HG2	1:67:A:GLY:HA3	13	0.25
(1,661)	1:66:A:MET:HG3	1:67:A:GLY:HA3	13	0.25
(1,340)	1:26:A:VAL:HG11	1:27:A:PRO:HD3	12	0.25
(1,340)	1:26:A:VAL:HG12	1:27:A:PRO:HD3	12	0.25
(1,340)	1:26:A:VAL:HG13	1:27:A:PRO:HD3	12	0.25
(1,340)	1:26:A:VAL:HG21	1:27:A:PRO:HD3	12	0.25
(1,340)	1:26:A:VAL:HG22	1:27:A:PRO:HD3	12	0.25
(1,340)	1:26:A:VAL:HG23	1:27:A:PRO:HD3	12	0.25
(1,175)	1:26:A:VAL:HG11	1:30:A:TRP:HD1	18	0.25
(1,175)	1:26:A:VAL:HG12	1:30:A:TRP:HD1	18	0.25
(1,175)	1:26:A:VAL:HG13	1:30:A:TRP:HD1	18	0.25
(1,175)	1:26:A:VAL:HG21	1:30:A:TRP:HD1	18	0.25
(1,175)	1:26:A:VAL:HG22	1:30:A:TRP:HD1	18	0.25
(1,175)	1:26:A:VAL:HG23	1:30:A:TRP:HD1	18	0.25
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD11	13	0.24
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD12	13	0.24
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD13	13	0.24
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD21	13	0.24
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD22	13	0.24
(1,564)	1:36:A:GLY:H	1:37:A:LEU:HD23	13	0.24
(1,661)	1:66:A:MET:HG2	1:67:A:GLY:HA3	5	0.23
(1,661)	1:66:A:MET:HG3	1:67:A:GLY:HA3	5	0.23
(1,661)	1:66:A:MET:HG2	1:67:A:GLY:HA3	17	0.23
(1,661)	1:66:A:MET:HG3	1:67:A:GLY:HA3	17	0.23
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG11	7	0.21
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG12	7	0.21
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG13	7	0.21
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG21	7	0.21
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG22	7	0.21
(1,456)	1:25:A:THR:HB	1:26:A:VAL:HG23	7	0.21
(1,175)	1:26:A:VAL:HG11	1:30:A:TRP:HD1	14	0.2
(1,175)	1:26:A:VAL:HG12	1:30:A:TRP:HD1	14	0.2
(1,175)	1:26:A:VAL:HG13	1:30:A:TRP:HD1	14	0.2
(1,175)	1:26:A:VAL:HG21	1:30:A:TRP:HD1	14	0.2
(1,175)	1:26:A:VAL:HG22	1:30:A:TRP:HD1	14	0.2
(1,175)	1:26:A:VAL:HG23	1:30:A:TRP:HD1	14	0.2
(1,178)	1:44:A:ARG:HA	1:44:A:ARG:HE	11	0.19

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,178)	1:44:A:ARG:HA	1:44:A:ARG:HE	13	0.19
(1,340)	1:26:A:VAL:HG11	1:27:A:PRO:HD3	11	0.16
(1,340)	1:26:A:VAL:HG12	1:27:A:PRO:HD3	11	0.16
(1,340)	1:26:A:VAL:HG13	1:27:A:PRO:HD3	11	0.16
(1,340)	1:26:A:VAL:HG21	1:27:A:PRO:HD3	11	0.16
(1,340)	1:26:A:VAL:HG22	1:27:A:PRO:HD3	11	0.16
(1,340)	1:26:A:VAL:HG23	1:27:A:PRO:HD3	11	0.16
(1,175)	1:26:A:VAL:HG11	1:30:A:TRP:HD1	5	0.16
(1,175)	1:26:A:VAL:HG12	1:30:A:TRP:HD1	5	0.16
(1,175)	1:26:A:VAL:HG13	1:30:A:TRP:HD1	5	0.16
(1,175)	1:26:A:VAL:HG21	1:30:A:TRP:HD1	5	0.16
(1,175)	1:26:A:VAL:HG22	1:30:A:TRP:HD1	5	0.16
(1,175)	1:26:A:VAL:HG23	1:30:A:TRP:HD1	5	0.16
(1,506)	1:23:A:GLU:HB2	1:24:A:CYS:H	18	0.15
(1,506)	1:23:A:GLU:HB3	1:24:A:CYS:H	18	0.15
(1,175)	1:26:A:VAL:HG11	1:30:A:TRP:HD1	6	0.15
(1,175)	1:26:A:VAL:HG12	1:30:A:TRP:HD1	6	0.15
(1,175)	1:26:A:VAL:HG13	1:30:A:TRP:HD1	6	0.15
(1,175)	1:26:A:VAL:HG21	1:30:A:TRP:HD1	6	0.15
(1,175)	1:26:A:VAL:HG22	1:30:A:TRP:HD1	6	0.15
(1,175)	1:26:A:VAL:HG23	1:30:A:TRP:HD1	6	0.15
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG2	4	0.15
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG3	4	0.15
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG2	5	0.15
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG3	5	0.15
(1,126)	1:30:A:TRP:H	1:30:A:TRP:HD1	18	0.15
(1,87)	1:23:A:GLU:H	1:23:A:GLU:HG2	15	0.15
(1,87)	1:23:A:GLU:H	1:23:A:GLU:HG3	15	0.15
(1,87)	1:23:A:GLU:H	1:23:A:GLU:HG2	18	0.15
(1,87)	1:23:A:GLU:H	1:23:A:GLU:HG3	18	0.15
(1,178)	1:44:A:ARG:HA	1:44:A:ARG:HE	7	0.14
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG2	2	0.14
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG3	2	0.14
(1,620)	1:51:A:CYS:HB2	1:52:A:CYS:H	10	0.13
(1,620)	1:51:A:CYS:HB3	1:52:A:CYS:H	10	0.13
(1,458)	1:26:A:VAL:HG11	1:30:A:TRP:HB2	19	0.13
(1,458)	1:26:A:VAL:HG12	1:30:A:TRP:HB2	19	0.13
(1,458)	1:26:A:VAL:HG13	1:30:A:TRP:HB2	19	0.13
(1,458)	1:26:A:VAL:HG21	1:30:A:TRP:HB2	19	0.13
(1,458)	1:26:A:VAL:HG22	1:30:A:TRP:HB2	19	0.13
(1,458)	1:26:A:VAL:HG23	1:30:A:TRP:HB2	19	0.13
(1,454)	1:26:A:VAL:HG11	1:30:A:TRP:HE3	19	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,454)	1:26:A:VAL:HG12	1:30:A:TRP:HE3	19	0.13
(1,454)	1:26:A:VAL:HG13	1:30:A:TRP:HE3	19	0.13
(1,454)	1:26:A:VAL:HG21	1:30:A:TRP:HE3	19	0.13
(1,454)	1:26:A:VAL:HG22	1:30:A:TRP:HE3	19	0.13
(1,454)	1:26:A:VAL:HG23	1:30:A:TRP:HE3	19	0.13
(1,340)	1:26:A:VAL:HG11	1:27:A:PRO:HD3	19	0.13
(1,340)	1:26:A:VAL:HG12	1:27:A:PRO:HD3	19	0.13
(1,340)	1:26:A:VAL:HG13	1:27:A:PRO:HD3	19	0.13
(1,340)	1:26:A:VAL:HG21	1:27:A:PRO:HD3	19	0.13
(1,340)	1:26:A:VAL:HG22	1:27:A:PRO:HD3	19	0.13
(1,340)	1:26:A:VAL:HG23	1:27:A:PRO:HD3	19	0.13
(1,661)	1:66:A:MET:HG2	1:67:A:GLY:HA3	18	0.12
(1,661)	1:66:A:MET:HG3	1:67:A:GLY:HA3	18	0.12
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD11	15	0.12
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD12	15	0.12
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD13	15	0.12
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD21	15	0.12
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD22	15	0.12
(1,563)	1:35:A:LYS:HE2	1:37:A:LEU:HD23	15	0.12
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD11	15	0.12
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD12	15	0.12
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD13	15	0.12
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD21	15	0.12
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD22	15	0.12
(1,563)	1:35:A:LYS:HE3	1:37:A:LEU:HD23	15	0.12
(1,454)	1:26:A:VAL:HG11	1:30:A:TRP:HE3	15	0.12
(1,454)	1:26:A:VAL:HG12	1:30:A:TRP:HE3	15	0.12
(1,454)	1:26:A:VAL:HG13	1:30:A:TRP:HE3	15	0.12
(1,454)	1:26:A:VAL:HG21	1:30:A:TRP:HE3	15	0.12
(1,454)	1:26:A:VAL:HG22	1:30:A:TRP:HE3	15	0.12
(1,454)	1:26:A:VAL:HG23	1:30:A:TRP:HE3	15	0.12
(1,208)	1:55:A:TYR:HD1	1:56:A:GLU:H	14	0.12
(1,208)	1:55:A:TYR:HD2	1:56:A:GLU:H	14	0.12
(1,629)	1:53:A:LYS:HG2	1:54:A:VAL:H	18	0.11
(1,629)	1:53:A:LYS:HG3	1:54:A:VAL:H	18	0.11
(1,581)	1:41:A:ARG:HB2	1:41:A:ARG:HE	16	0.11
(1,581)	1:41:A:ARG:HB3	1:41:A:ARG:HE	16	0.11
(1,458)	1:26:A:VAL:HG11	1:30:A:TRP:HB2	5	0.11
(1,458)	1:26:A:VAL:HG12	1:30:A:TRP:HB2	5	0.11
(1,458)	1:26:A:VAL:HG13	1:30:A:TRP:HB2	5	0.11
(1,458)	1:26:A:VAL:HG21	1:30:A:TRP:HB2	5	0.11
(1,458)	1:26:A:VAL:HG22	1:30:A:TRP:HB2	5	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,458)	1:26:A:VAL:HG23	1:30:A:TRP:HB2	5	0.11
(1,358)	1:46:A:TYR:HA	1:63:A:TYR:HB2	14	0.11
(1,340)	1:26:A:VAL:HG11	1:27:A:PRO:HD3	1	0.11
(1,340)	1:26:A:VAL:HG12	1:27:A:PRO:HD3	1	0.11
(1,340)	1:26:A:VAL:HG13	1:27:A:PRO:HD3	1	0.11
(1,340)	1:26:A:VAL:HG21	1:27:A:PRO:HD3	1	0.11
(1,340)	1:26:A:VAL:HG22	1:27:A:PRO:HD3	1	0.11
(1,340)	1:26:A:VAL:HG23	1:27:A:PRO:HD3	1	0.11
(1,279)	1:65:A:ARG:HG2	1:68:A:GLU:H	18	0.11
(1,279)	1:65:A:ARG:HG3	1:68:A:GLU:H	18	0.11
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG2	6	0.11
(1,165)	1:65:A:ARG:H	1:65:A:ARG:HG3	6	0.11
(1,85)	1:25:A:THR:HB	1:26:A:VAL:H	17	0.11
(1,64)	1:23:A:GLU:HG2	1:24:A:CYS:H	11	0.11
(1,64)	1:23:A:GLU:HG3	1:24:A:CYS:H	11	0.11
(1,408)	1:35:A:LYS:HD2	1:38:A:CYS:H	5	0.1
(1,408)	1:35:A:LYS:HD3	1:38:A:CYS:H	5	0.1
(1,392)	1:65:A:ARG:HB2	1:67:A:GLY:H	4	0.1
(1,341)	1:47:A:CYS:H	1:47:A:CYS:HB3	13	0.1
(1,327)	1:44:A:ARG:HA	1:44:A:ARG:HD2	15	0.1
(1,311)	1:58:A:CYS:H	1:58:A:CYS:HB3	16	0.1
(1,227)	1:45:A:TYR:HD1	1:65:A:ARG:HG2	20	0.1
(1,227)	1:45:A:TYR:HD1	1:65:A:ARG:HG3	20	0.1
(1,227)	1:45:A:TYR:HD2	1:65:A:ARG:HG2	20	0.1
(1,227)	1:45:A:TYR:HD2	1:65:A:ARG:HG3	20	0.1

10 Dihedral-angle violation analysis (i)

No dihedral-angle restraints found