



# Full wwPDB NMR Structure Validation Report ⓘ

Dec 24, 2024 – 02:09 PM EST

PDB ID : 2KZH  
BMRB ID : 17005  
Title : Three-dimensional structure of a truncated phosphoribosylanthranilate isomerase (residues 255-384) from Escherichia coli  
Authors : Setiyaputra, S.; Mackay, J.P.; Patrick, W.M.  
Deposited on : 2010-06-17

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
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) : 2.40

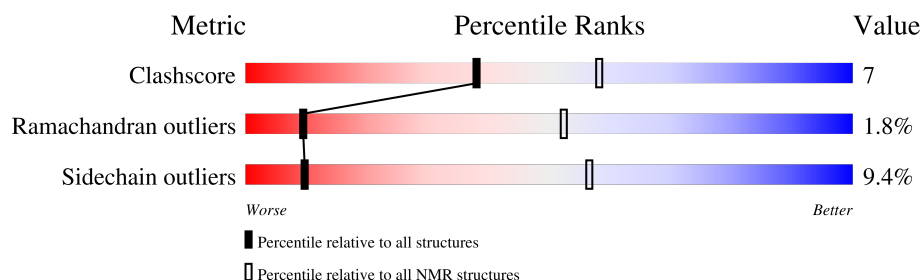
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 85%.

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	210492	14027
Ramachandran outliers	207382	12486
Sidechain outliers	206894	12463

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	140	

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 13 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:4-A:106, A:117-A:125 (112)	0.75	13

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 3 clusters. No single-model clusters were found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1999 atoms, of which 994 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Tryptophan biosynthesis protein trpCF.

Mol	Chain	Residues	Atoms						Trace
1	A	134	Total	C	H	N	O	S	0
			1999	632	994	177	192	4	

There are 10 discrepancies between the modelled and reference sequences:

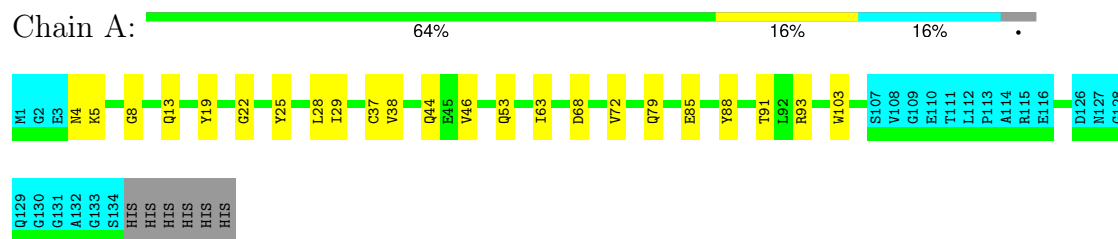
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P00909
A	132	ALA	-	SEE REMARK 999	UNP P00909
A	133	GLY	-	SEE REMARK 999	UNP P00909
A	134	SER	-	SEE REMARK 999	UNP P00909
A	135	HIS	-	expression tag	UNP P00909
A	136	HIS	-	expression tag	UNP P00909
A	137	HIS	-	expression tag	UNP P00909
A	138	HIS	-	expression tag	UNP P00909
A	139	HIS	-	expression tag	UNP P00909
A	140	HIS	-	expression tag	UNP P00909

## 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: Tryptophan biosynthesis protein trpCF

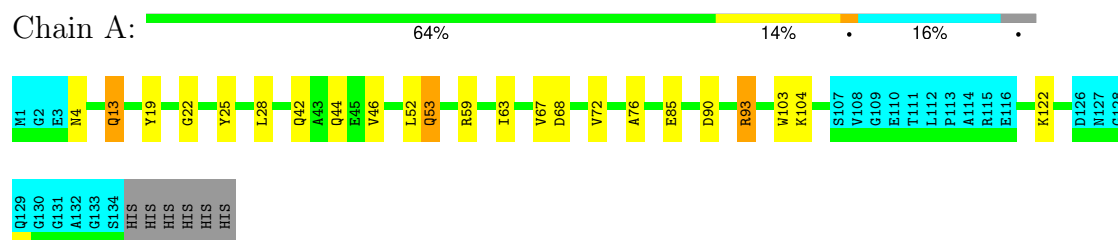


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

#### 4.2.1 Score per residue for model 1

- Molecule 1: Tryptophan biosynthesis protein trpCF



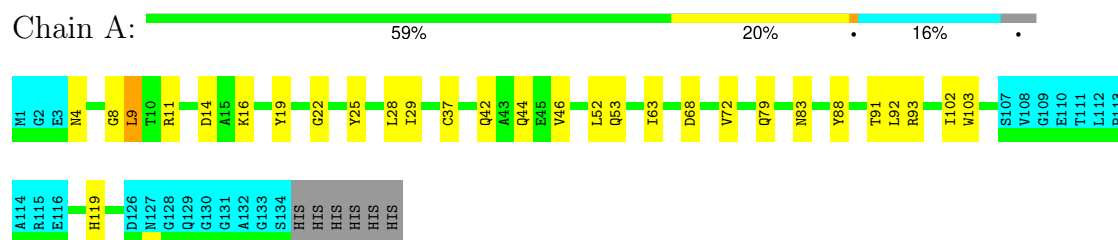
#### 4.2.2 Score per residue for model 2

- Molecule 1: Tryptophan biosynthesis protein trpCF



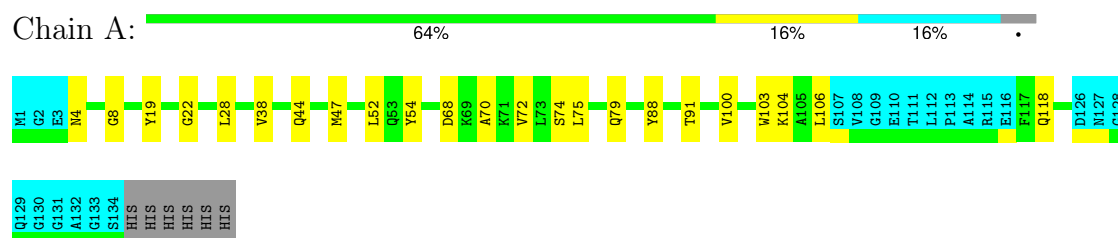
### 4.2.3 Score per residue for model 3

- Molecule 1: Tryptophan biosynthesis protein trpCF



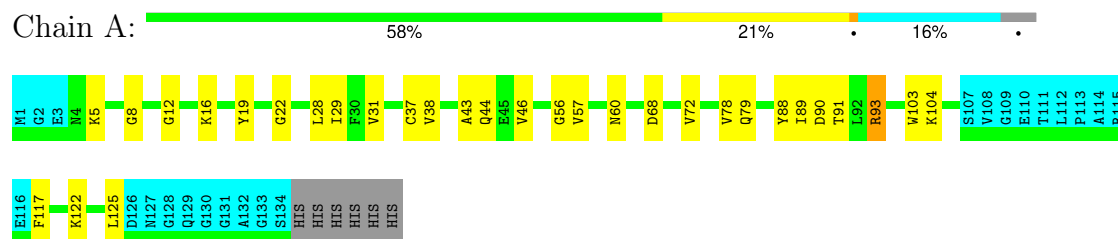
#### 4.2.4 Score per residue for model 4

- Molecule 1: Tryptophan biosynthesis protein trpCF



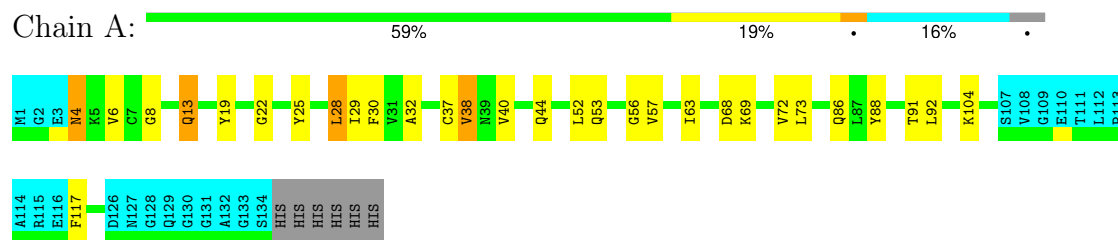
#### 4.2.5 Score per residue for model 5

- Molecule 1: Tryptophan biosynthesis protein trpCF



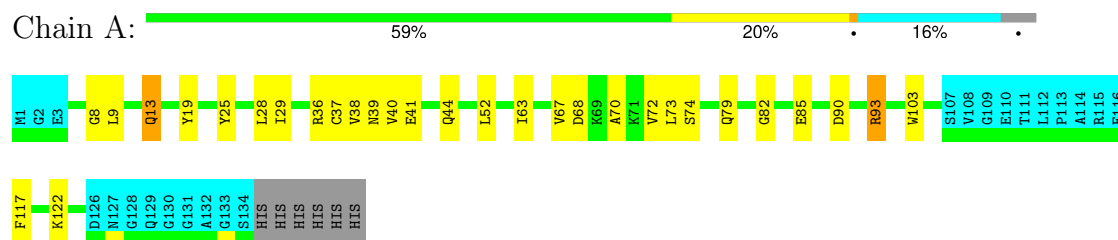
#### 4.2.6 Score per residue for model 6

- Molecule 1: Tryptophan biosynthesis protein trpCF



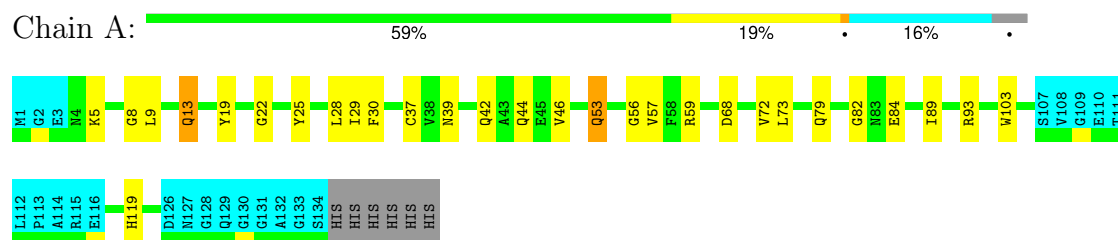
#### 4.2.7 Score per residue for model 7

- Molecule 1: Tryptophan biosynthesis protein trpCF



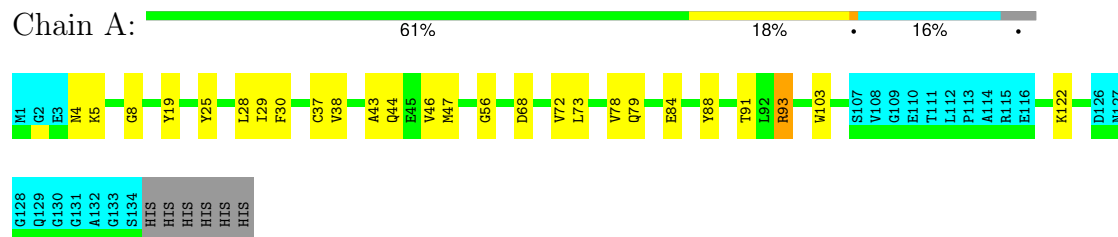
#### 4.2.8 Score per residue for model 8

- Molecule 1: Tryptophan biosynthesis protein trpCF



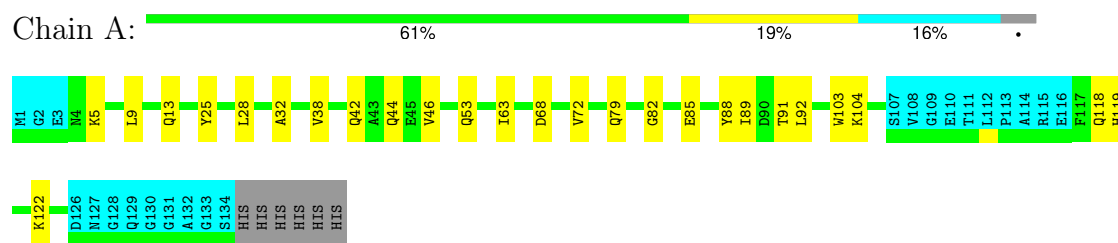
#### 4.2.9 Score per residue for model 9

- Molecule 1: Tryptophan biosynthesis protein trpCF



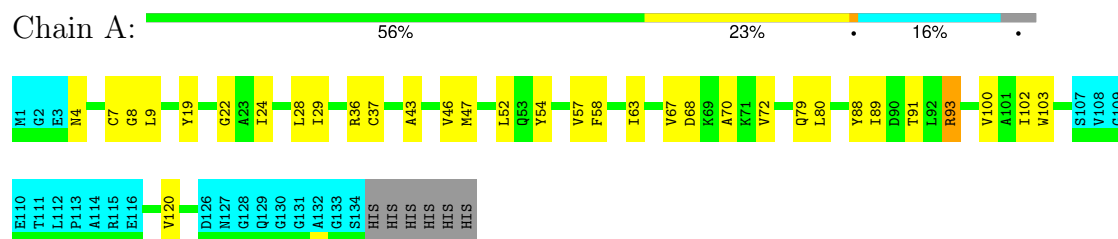
### 4.2.10 Score per residue for model 10

- Molecule 1: Tryptophan biosynthesis protein trpCF



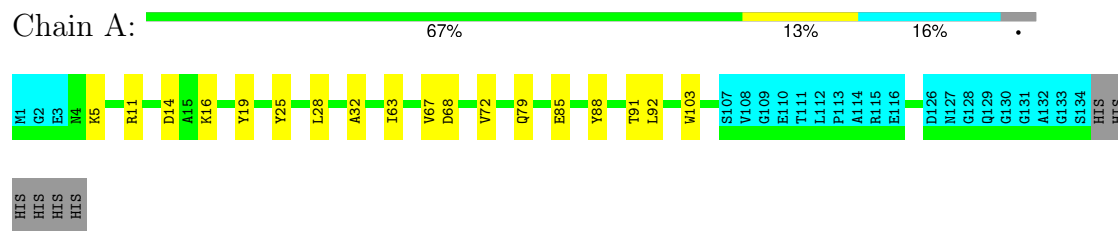
### 4.2.11 Score per residue for model 11

- Molecule 1: Tryptophan biosynthesis protein trpCF



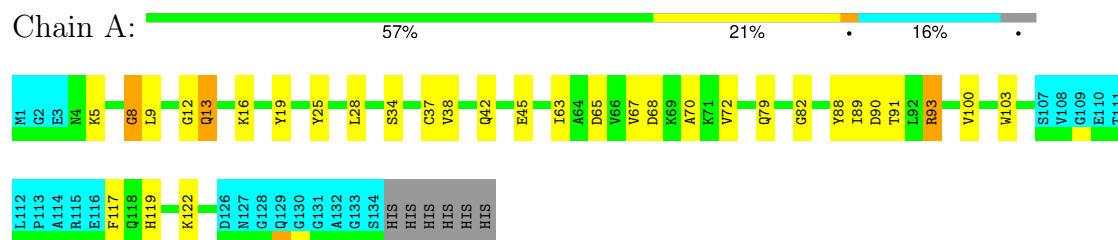
### 4.2.12 Score per residue for model 12

- Molecule 1: Tryptophan biosynthesis protein trpCF



### 4.2.13 Score per residue for model 13 (medoid)

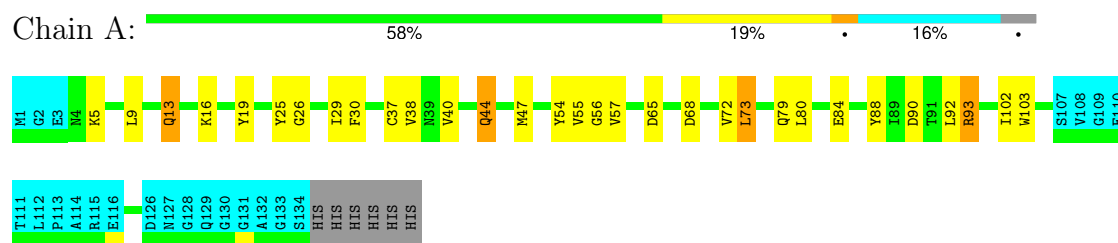
- Molecule 1: Tryptophan biosynthesis protein trpCF





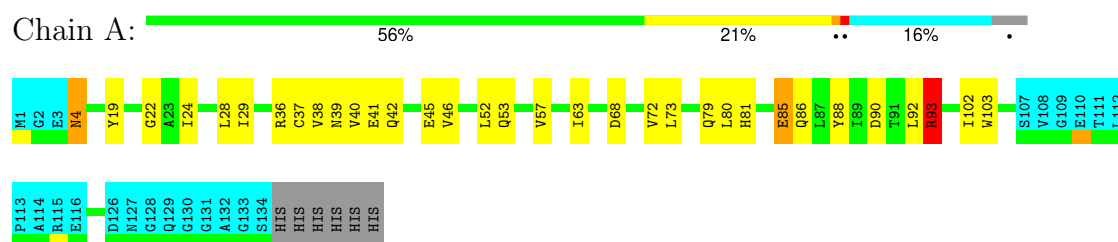
#### 4.2.14 Score per residue for model 14

- Molecule 1: Tryptophan biosynthesis protein trpCF



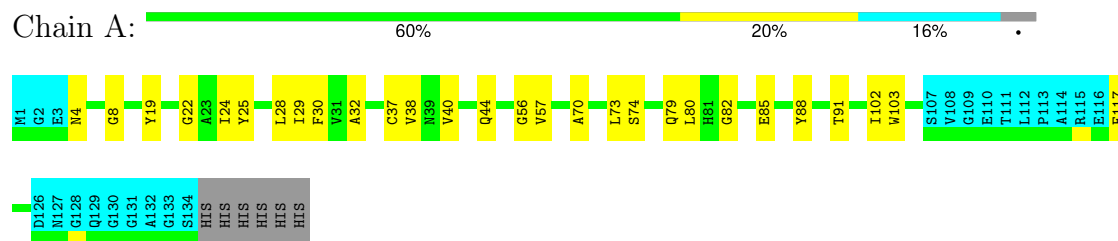
#### 4.2.15 Score per residue for model 15

- Molecule 1: Tryptophan biosynthesis protein trpCF



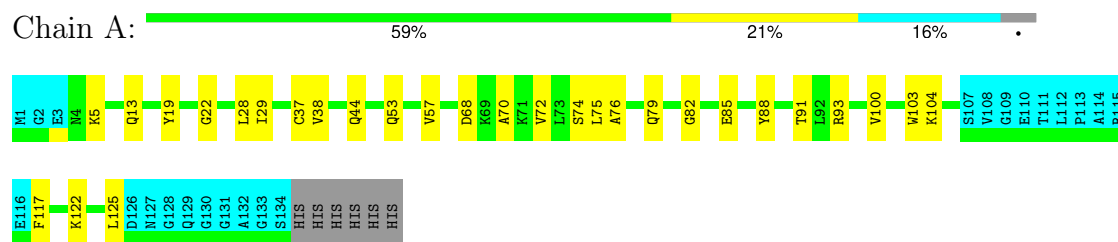
#### 4.2.16 Score per residue for model 16

- Molecule 1: Tryptophan biosynthesis protein trpCF



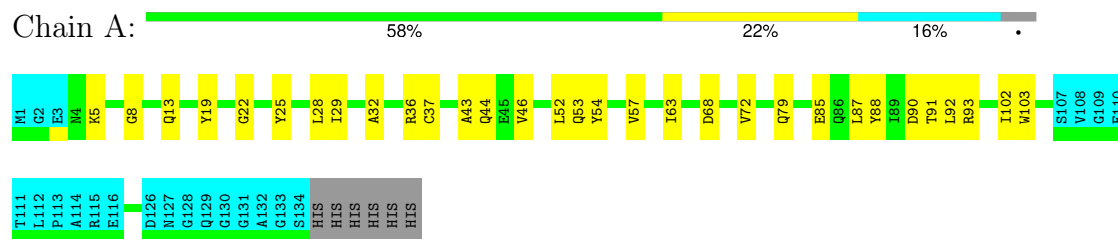
#### 4.2.17 Score per residue for model 17

- Molecule 1: Tryptophan biosynthesis protein trpCF



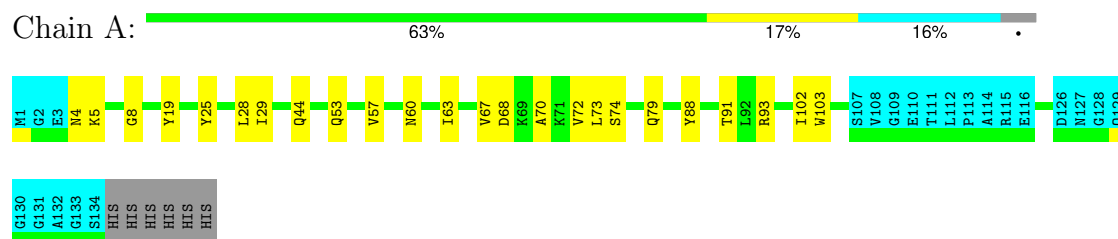
### 4.2.18 Score per residue for model 18

- Molecule 1: Tryptophan biosynthesis protein trpCF



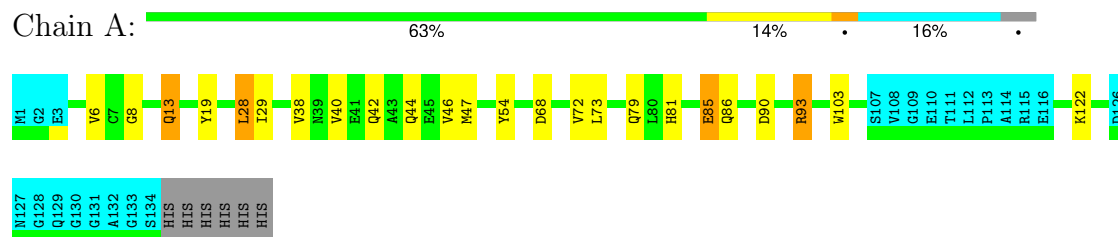
### 4.2.19 Score per residue for model 19

- Molecule 1: Tryptophan biosynthesis protein trpCF



### 4.2.20 Score per residue for model 20

- Molecule 1: Tryptophan biosynthesis protein trpCF



## 5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *structures with the least restraint violations*.

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

Software name	Classification	Version
CNS	structure solution	1.21
CNS	refinement	1.21

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	1482
Number of shifts mapped to atoms	1482
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	85%

## 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	0.91±0.02	0±0/876 ( 0.0± 0.0%)	0.74±0.03	0±0/1193 ( 0.0± 0.0%)
All	All	0.91	1/17520 ( 0.0%)	0.74	0/23860 ( 0.0%)

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

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	0.1±0.4
All	All	0	3

All unique bond outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	39	ASN	CG-ND2	-5.18	1.20	1.32	8	1

There are no bond-angle outliers.

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	93	ARG	Sidechain	2
1	A	59	ARG	Sidechain	1

### 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	859	862	862	12±4
All	All	17180	17240	17240	247

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 7.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:43:ALA:HA	1:A:46:VAL:HG12	0.75	1.57	2	5
1:A:93:ARG:NH2	1:A:102:ILE:HB	0.69	2.02	11	1
1:A:76:ALA:O	1:A:100:VAL:HA	0.66	1.91	17	1
1:A:47:MET:HA	1:A:54:TYR:OH	0.63	1.93	11	3
1:A:90:ASP:HA	1:A:93:ARG:NE	0.62	2.09	5	3
1:A:93:ARG:HD2	1:A:102:ILE:HD12	0.62	1.72	3	2
1:A:13:GLN:H	1:A:13:GLN:NE2	0.61	1.94	2	10
1:A:47:MET:SD	1:A:73:LEU:HA	0.59	2.37	9	1
1:A:90:ASP:HA	1:A:93:ARG:NH1	0.57	2.15	15	2
1:A:90:ASP:HA	1:A:93:ARG:CZ	0.57	2.29	14	5
1:A:79:GLN:NE2	1:A:103:TRP:HB2	0.56	2.15	14	18
1:A:42:GLN:O	1:A:45:GLU:HG2	0.55	2.01	13	2
1:A:82:GLY:HA3	1:A:117:PHE:CZ	0.55	2.37	7	4
1:A:42:GLN:O	1:A:46:VAL:HG23	0.54	2.02	1	6
1:A:53:GLN:NE2	1:A:76:ALA:HB2	0.54	2.16	1	1
1:A:29:ILE:O	1:A:37:CYS:HB3	0.54	2.03	16	10
1:A:79:GLN:HE22	1:A:103:TRP:HB2	0.53	1.63	20	3
1:A:93:ARG:NH2	1:A:120:VAL:HA	0.53	2.18	11	1
1:A:93:ARG:HD2	1:A:102:ILE:HD13	0.53	1.79	18	2
1:A:63:ILE:HG13	1:A:92:LEU:HD13	0.53	1.81	3	7
1:A:70:ALA:O	1:A:74:SER:HA	0.53	2.03	7	5
1:A:40:VAL:HG23	1:A:73:LEU:HB2	0.52	1.81	7	4
1:A:70:ALA:HB1	1:A:100:VAL:HG21	0.52	1.79	13	5
1:A:4:ASN:HD21	1:A:24:ILE:HB	0.52	1.65	15	1
1:A:29:ILE:HG23	1:A:57:VAL:HB	0.52	1.82	6	2
1:A:30:PHE:CE2	1:A:56:GLY:HA3	0.51	2.40	16	5
1:A:4:ASN:OD1	1:A:24:ILE:HB	0.51	2.06	16	2
1:A:4:ASN:HD22	1:A:4:ASN:N	0.50	2.04	15	2
1:A:40:VAL:HG21	1:A:69:LYS:HD2	0.50	1.83	6	1
1:A:29:ILE:HG22	1:A:57:VAL:O	0.50	2.05	8	4
1:A:88:TYR:O	1:A:91:THR:HB	0.50	2.07	10	14

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:68:ASP:O	1:A:72:VAL:HG23	0.50	2.06	12	19
1:A:12:GLY:O	1:A:16:LYS:HG3	0.49	2.07	13	3
1:A:29:ILE:CG2	1:A:57:VAL:HB	0.49	2.38	14	6
1:A:103:TRP:CD1	1:A:122:LYS:HB2	0.48	2.43	13	5
1:A:82:GLY:HA3	1:A:117:PHE:CE1	0.47	2.45	2	1
1:A:28:LEU:HD11	1:A:54:TYR:CD2	0.47	2.45	20	1
1:A:88:TYR:CZ	1:A:92:LEU:HD21	0.47	2.45	15	2
1:A:103:TRP:CD1	1:A:122:LYS:HB3	0.46	2.45	5	4
1:A:25:TYR:HA	1:A:53:GLN:HB3	0.46	1.87	8	1
1:A:63:ILE:CD1	1:A:91:THR:HG22	0.46	2.41	3	2
1:A:34:SER:OG	1:A:37:CYS:HB2	0.46	2.10	13	1
1:A:6:VAL:O	1:A:28:LEU:HB3	0.46	2.11	20	2
1:A:82:GLY:HA3	1:A:117:PHE:CE2	0.46	2.46	13	1
1:A:8:GLY:HA2	1:A:38:VAL:HG23	0.46	1.88	6	2
1:A:36:ARG:HA	1:A:36:ARG:HE	0.46	1.70	18	1
1:A:36:ARG:HG3	1:A:37:CYS:SG	0.45	2.51	15	2
1:A:93:ARG:CZ	1:A:119:HIS:HB2	0.45	2.42	13	1
1:A:82:GLY:HA2	1:A:89:ILE:HD11	0.45	1.87	13	3
1:A:118:GLN:HG3	1:A:119:HIS:CD2	0.45	2.47	10	1
1:A:63:ILE:O	1:A:67:VAL:HG23	0.45	2.11	7	7
1:A:65:ASP:O	1:A:68:ASP:HB3	0.45	2.12	13	2
1:A:79:GLN:HA	1:A:103:TRP:O	0.44	2.12	4	2
1:A:44:GLN:O	1:A:47:MET:HG2	0.44	2.12	14	1
1:A:80:LEU:HD13	1:A:92:LEU:HG	0.44	1.88	14	2
1:A:39:ASN:ND2	1:A:41:GLU:HB2	0.44	2.27	7	2
1:A:40:VAL:HG23	1:A:73:LEU:HD23	0.44	1.89	14	1
1:A:5:LYS:HG3	1:A:27:GLY:HA3	0.44	1.90	2	1
1:A:11:ARG:HB2	1:A:14:ASP:HB3	0.44	1.90	12	2
1:A:43:ALA:HA	1:A:46:VAL:CG1	0.43	2.38	2	2
1:A:8:GLY:HA3	1:A:37:CYS:N	0.43	2.28	18	1
1:A:85:GLU:CD	1:A:86:GLN:H	0.43	2.16	15	2
1:A:93:ARG:HG3	1:A:102:ILE:HD12	0.43	1.89	15	1
1:A:93:ARG:NH2	1:A:119:HIS:HB2	0.43	2.28	13	1
1:A:26:GLY:O	1:A:55:VAL:HG12	0.43	2.13	14	1
1:A:93:ARG:CD	1:A:102:ILE:HD12	0.43	2.42	2	2
1:A:9:LEU:HD11	1:A:14:ASP:HB3	0.43	1.90	3	1
1:A:29:ILE:HG13	1:A:37:CYS:SG	0.43	2.53	17	2
1:A:80:LEU:CD1	1:A:92:LEU:HG	0.43	2.44	15	1
1:A:4:ASN:ND2	1:A:24:ILE:HB	0.42	2.27	15	1
1:A:85:GLU:HB3	1:A:88:TYR:HB2	0.42	1.92	16	1
1:A:56:GLY:O	1:A:78:VAL:HA	0.42	2.15	5	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:93:ARG:NH1	1:A:119:HIS:HB2	0.42	2.30	8	1
1:A:87:LEU:O	1:A:90:ASP:HB3	0.42	2.15	18	1
1:A:59:ARG:HD3	1:A:84:GLU:OE1	0.41	2.15	8	1
1:A:11:ARG:HB2	1:A:14:ASP:CB	0.41	2.46	3	1
1:A:80:LEU:HD12	1:A:102:ILE:HG12	0.41	1.92	14	1
1:A:93:ARG:CZ	1:A:102:ILE:HB	0.41	2.45	11	1
1:A:93:ARG:H	1:A:93:ARG:HD3	0.41	1.75	9	1
1:A:40:VAL:CG2	1:A:69:LYS:HD2	0.41	2.46	6	1
1:A:86:GLN:HA	1:A:117:PHE:CD1	0.41	2.50	6	1
1:A:7:CYS:HB2	1:A:36:ARG:O	0.41	2.15	11	1
1:A:80:LEU:CD2	1:A:89:ILE:HG13	0.41	2.46	11	1
1:A:5:LYS:HD3	1:A:25:TYR:OH	0.40	2.16	2	1
1:A:44:GLN:HE21	1:A:44:GLN:HB3	0.40	1.60	14	1
1:A:80:LEU:HD23	1:A:102:ILE:HG23	0.40	1.93	16	1
1:A:89:ILE:HD13	1:A:117:PHE:CD1	0.40	2.52	5	1
1:A:8:GLY:O	1:A:9:LEU:HB3	0.40	2.15	13	1
1:A:79:GLN:OE1	1:A:103:TRP:HB2	0.40	2.16	15	1
1:A:31:VAL:CG1	1:A:37:CYS:HB2	0.40	2.47	5	1
1:A:93:ARG:HG3	1:A:119:HIS:CE1	0.40	2.51	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	112/140 (80%)	100±2 (89±1%)	10±2 (9±1%)	2±1 (2±1%)	9	52
All	All	2240/2800 (80%)	1990 (89%)	210 (9%)	40 (2%)	9	52

All 6 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	22	GLY	11
1	A	8	GLY	11
1	A	5	LYS	10

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Mol	Chain	Res	Type	Models (Total)
1	A	32	ALA	5
1	A	81	HIS	2
1	A	83	ASN	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	87/107 (81%)	79±2 (91±2%)	8±2 (9±2%)	10	56
All	All	1740/2140 (81%)	1577 (91%)	163 (9%)	10	56

All 25 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	19	TYR	19
1	A	28	LEU	19
1	A	44	GLN	15
1	A	25	TYR	13
1	A	38	VAL	11
1	A	13	GLN	9
1	A	53	GLN	9
1	A	85	GLU	9
1	A	93	ARG	9
1	A	52	LEU	8
1	A	4	ASN	7
1	A	9	LEU	7
1	A	104	LYS	6
1	A	73	LEU	4
1	A	16	LYS	3
1	A	47	MET	2
1	A	75	LEU	2
1	A	60	ASN	2
1	A	125	LEU	2
1	A	84	GLU	2
1	A	106	LEU	1
1	A	118	GLN	1

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Mol	Chain	Res	Type	Models (Total)
1	A	58	PHE	1
1	A	54	TYR	1
1	A	29	ILE	1

### 6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 6.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 6.7 Other polymers [i](#)

There are no such molecules in this entry.

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

There are no chain breaks in this entry.

## 7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 85% for the well-defined parts and 85% for the entire structure.

### 7.1 Chemical shift list 1

File name: working\_cs.cif

Chemical shift list name: *assigned\_chem\_shift\_list\_1*

#### 7.1.1 Bookkeeping

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

Total number of shifts	1482
Number of shifts mapped to atoms	1482
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
$^{13}\text{C}_\alpha$	132	$-0.27 \pm 0.08$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}_\beta$	119	$0.06 \pm 0.07$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}'$	134	$-0.08 \pm 0.13$	None needed ( $< 0.5$ ppm)
$^{15}\text{N}$	125	$-0.05 \pm 0.38$	None needed ( $< 0.5$ ppm)

#### 7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 85%, i.e. 1289 atoms were assigned a chemical shift out of a possible 1513. 0 out of 25 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	552/561 (98%)	224/228 (98%)	223/224 (100%)	105/109 (96%)
Sidechain	699/837 (84%)	476/551 (86%)	214/257 (83%)	9/29 (31%)

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	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Aromatic	38/115 (33%)	37/57 (65%)	0/53 (0%)	1/5 (20%)
Overall	1289/1513 (85%)	737/836 (88%)	437/534 (82%)	115/143 (80%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 85%, i.e. 1482 atoms were assigned a chemical shift out of a possible 1749. 0 out of 27 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	661/675 (98%)	270/277 (97%)	266/268 (99%)	125/130 (96%)
Sidechain	783/959 (82%)	535/629 (85%)	237/296 (80%)	11/34 (32%)
Aromatic	38/115 (33%)	37/57 (65%)	0/53 (0%)	1/5 (20%)
Overall	1482/1749 (85%)	842/963 (87%)	503/617 (82%)	137/169 (81%)

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

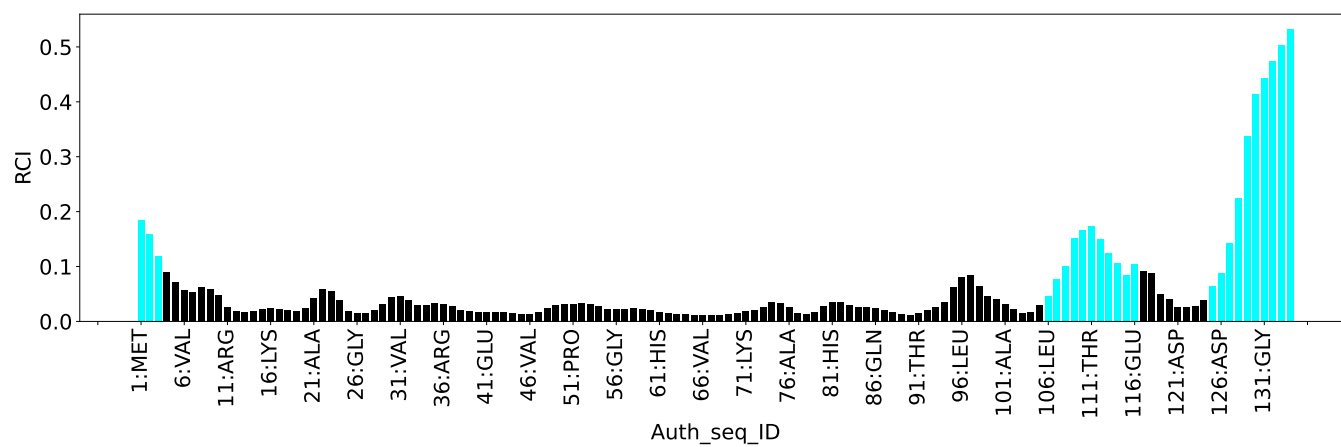
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	122	LYS	HB2	0.39	0.58 – 2.97	-5.8
1	A	77	ALA	HB1	0.04	0.14 – 2.58	-5.4
1	A	77	ALA	HB2	0.04	0.14 – 2.58	-5.4
1	A	77	ALA	HB3	0.04	0.14 – 2.58	-5.4
1	A	122	LYS	HD2	0.55	0.58 – 2.64	-5.1
1	A	115	ARG	HG2	0.23	0.26 – 2.87	-5.1

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

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

Random coil index (RCI) for chain A:



## 8 NMR restraints analysis

### 8.1 Conformationally restricting restraints

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

Description	Value
Total distance restraints	2297
Intra-residue ( $ i-j =0$ )	515
Sequential ( $ i-j =1$ )	595
Medium range ( $ i-j >1$ and $ i-j <5$ )	527
Long range ( $ i-j \geq 5$ )	631
Inter-chain	0
Hydrogen bond restraints	29
Disulfide bond restraints	0
Total dihedral-angle restraints	208
Number of unmapped restraints	0
Number of restraints per residue	17.9
Number of long range restraints per residue <sup>1</sup>	4.6

<sup>1</sup>Long range hydrogen bonds and disulfide bonds are counted as long range restraints while calculating the number of long range restraints per residue

### 8.2 Residual restraint violations

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

#### 8.2.1 Average number of distance violations per model

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

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	13.1	0.2
0.2-0.5 (Medium)	14.9	0.5
>0.5 (Large)	12.2	3.64

### 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)	10.0	4.44
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

## 9 Distance violation analysis ⓘ

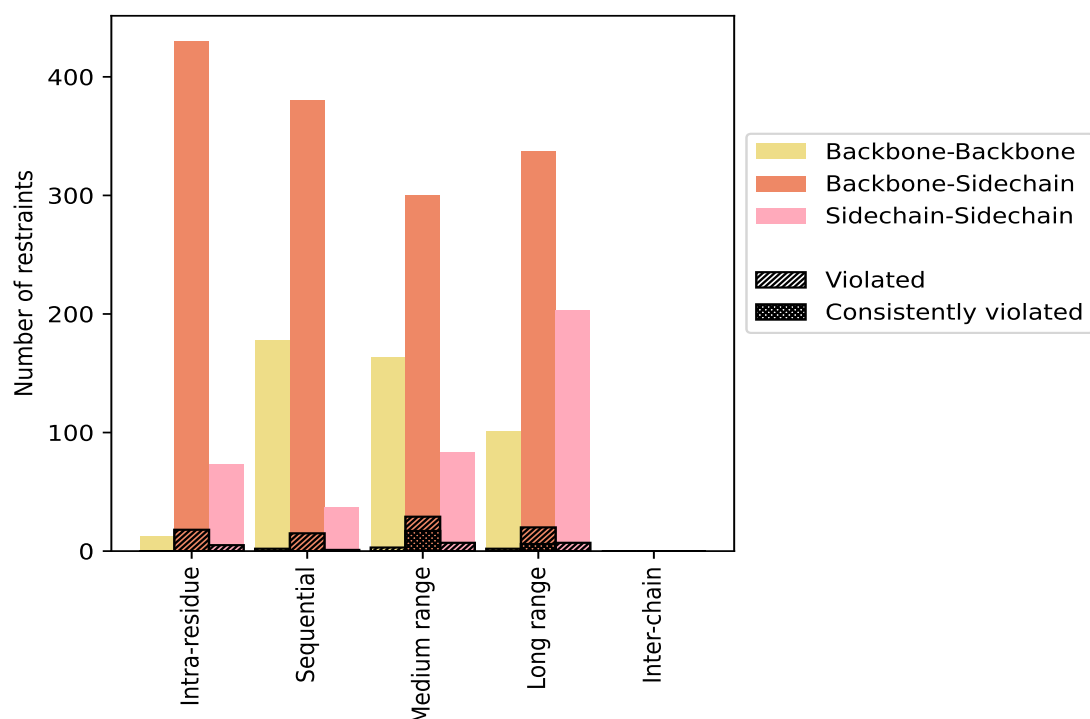
### 9.1 Summary of distance violations ⓘ

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

Restrains type	Count	% <sup>1</sup>	Violated <sup>3</sup>			Consistently Violated <sup>4</sup>		
			Count	% <sup>2</sup>	% <sup>1</sup>	Count	% <sup>2</sup>	% <sup>1</sup>
<a href="#">Intra-residue ( i-j =0)</a>	<a href="#">515</a>	<a href="#">22.4</a>	<a href="#">23</a>	<a href="#">4.5</a>	<a href="#">1.0</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0.0</a>
Backbone-Backbone	12	0.5	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	430	18.7	18	4.2	0.8	0	0.0	0.0
Sidechain-Sidechain	73	3.2	5	6.8	0.2	0	0.0	0.0
<a href="#">Sequential ( i-j =1)</a>	<a href="#">595</a>	<a href="#">25.9</a>	<a href="#">18</a>	<a href="#">3.0</a>	<a href="#">0.8</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0.0</a>
Backbone-Backbone	178	7.7	2	1.1	0.1	0	0.0	0.0
Backbone-Sidechain	380	16.5	15	3.9	0.7	0	0.0	0.0
Sidechain-Sidechain	37	1.6	1	2.7	0.0	0	0.0	0.0
<a href="#">Medium range ( i-j &gt;1 &amp;  i-j &lt;5)</a>	<a href="#">527</a>	<a href="#">22.9</a>	<a href="#">20</a>	<a href="#">3.8</a>	<a href="#">0.9</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0.0</a>
Backbone-Backbone	163	7.1	3	1.8	0.1	0	0.0	0.0
Backbone-Sidechain	281	12.2	10	3.6	0.4	0	0.0	0.0
Sidechain-Sidechain	83	3.6	7	8.4	0.3	0	0.0	0.0
<a href="#">Long range ( i-j ≥5)</a>	<a href="#">631</a>	<a href="#">27.5</a>	<a href="#">22</a>	<a href="#">3.5</a>	<a href="#">1.0</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0.0</a>
Backbone-Backbone	101	4.4	2	2.0	0.1	0	0.0	0.0
Backbone-Sidechain	327	14.2	13	4.0	0.6	0	0.0	0.0
Sidechain-Sidechain	203	8.8	7	3.4	0.3	0	0.0	0.0
<a href="#">Inter-chain</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0.0</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0.0</a>
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
<a href="#">Hydrogen bond</a>	<a href="#">29</a>	<a href="#">1.3</a>	<a href="#">26</a>	<a href="#">89.7</a>	<a href="#">1.1</a>	<a href="#">23</a>	<a href="#">79.3</a>	<a href="#">1.0</a>
<a href="#">Disulfide bond</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0.0</a>	<a href="#">0</a>	<a href="#">0.0</a>	<a href="#">0.0</a>
<a href="#">Total</a>	<a href="#">2297</a>	<a href="#">100.0</a>	<a href="#">109</a>	<a href="#">4.7</a>	<a href="#">4.7</a>	<a href="#">23</a>	<a href="#">1.0</a>	<a href="#">1.0</a>
Backbone-Backbone	454	19.8	7	1.5	0.3	0	0.0	0.0
Backbone-Sidechain	1447	63.0	82	5.7	3.6	23	1.6	1.0
Sidechain-Sidechain	396	17.2	20	5.1	0.9	0	0.0	0.0

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

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



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and 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 <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total				
1	3	6	19	13	0	41	0.45	1.62	0.39	0.31
2	4	4	19	12	0	39	0.47	2.19	0.45	0.35
3	3	7	20	14	0	44	0.46	1.77	0.44	0.3
4	8	5	18	10	0	41	0.43	1.79	0.39	0.28
5	7	6	21	10	0	44	0.39	1.59	0.36	0.25
6	2	7	19	14	0	42	0.42	1.75	0.39	0.27
7	5	5	19	15	0	44	0.45	1.62	0.37	0.32
8	6	4	24	13	0	47	0.39	1.45	0.33	0.27
9	4	2	21	9	0	36	0.55	1.95	0.47	0.43
10	3	9	20	14	0	46	0.44	1.85	0.44	0.25

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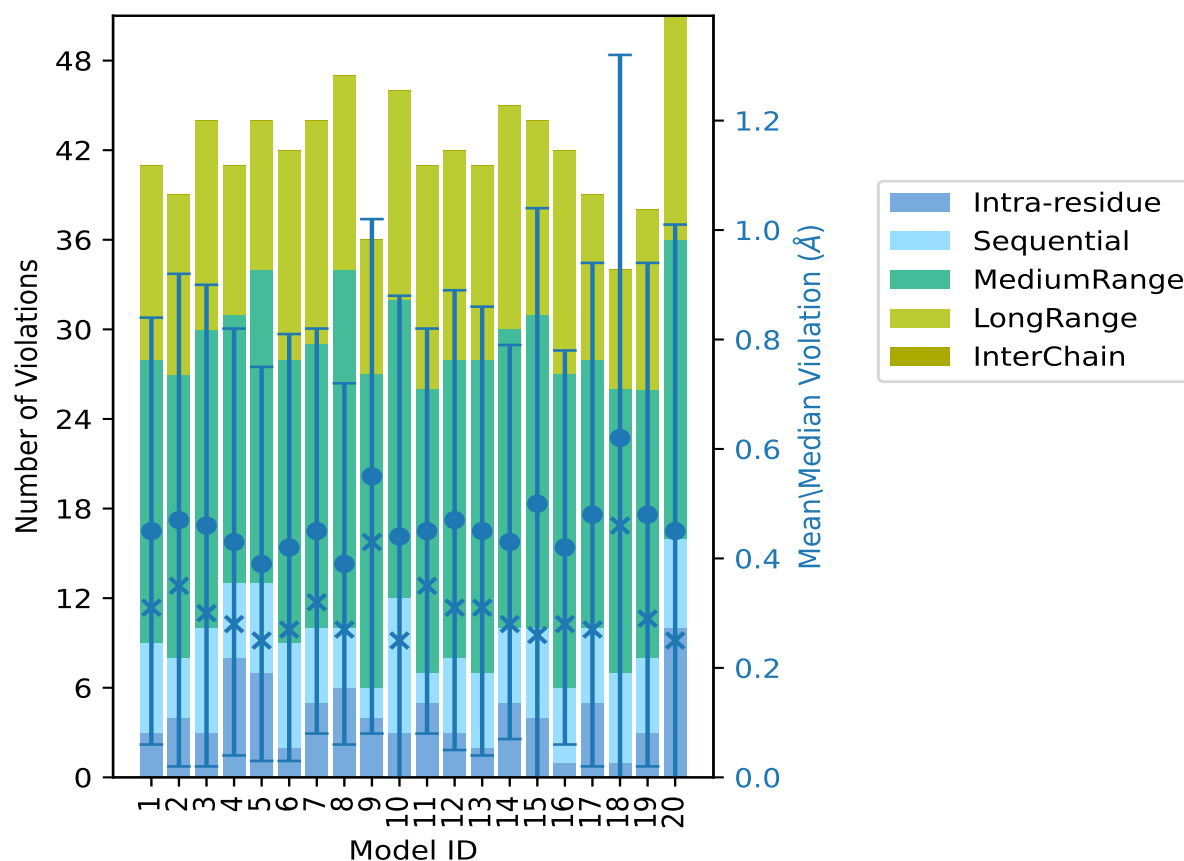
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total				
11	5	2	19	15	0	41	0.45	1.45	0.37	0.35
12	3	5	20	14	0	42	0.47	2.02	0.42	0.31
13	2	5	21	13	0	41	0.45	2.01	0.41	0.31
14	5	5	20	15	0	45	0.43	1.3	0.36	0.28
15	4	6	21	13	0	44	0.5	2.79	0.54	0.26
16	1	5	21	15	0	42	0.42	1.66	0.36	0.28
17	5	5	18	11	0	39	0.48	1.87	0.46	0.27
18	1	6	19	8	0	34	0.62	3.64	0.7	0.46
19	3	5	18	12	0	38	0.48	1.82	0.46	0.29
20	10	6	20	15	0	51	0.45	3.36	0.56	0.25

<sup>1</sup>Intra-residue restraints, <sup>2</sup>Sequential restraints, <sup>3</sup>Medium range restraints, <sup>4</sup>Long range restraints,

<sup>5</sup>Inter-chain restraints, <sup>6</sup>Standard deviation

### 9.2.1 Bar graph : Distance Violation statistics for each model ⓘ



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

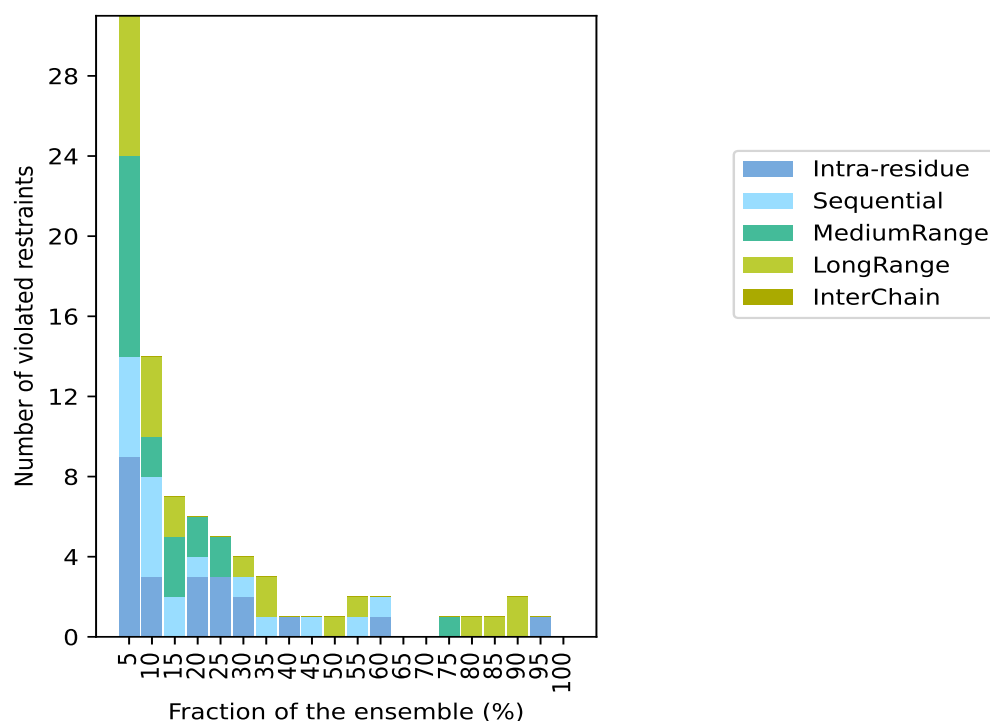
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 2185(IR:492, SQ:577, MR:507, LR:609, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total	Count <sup>6</sup>	%
9	5	10	7	0	31	1	5.0
3	5	2	4	0	14	2	10.0
0	2	3	2	0	7	3	15.0
3	1	2	0	0	6	4	20.0
3	0	2	0	0	5	5	25.0
2	1	0	1	0	4	6	30.0
0	1	0	2	0	3	7	35.0
1	0	0	0	0	1	8	40.0
0	1	0	0	0	1	9	45.0
0	0	0	1	0	1	10	50.0
0	1	0	1	0	2	11	55.0
1	1	0	0	0	2	12	60.0
0	0	0	0	0	0	13	65.0
0	0	0	0	0	0	14	70.0
0	0	1	0	0	1	15	75.0
0	0	0	1	0	1	16	80.0
0	0	0	1	0	1	17	85.0
0	0	0	2	0	2	18	90.0
1	0	0	0	0	1	19	95.0
0	0	0	0	0	0	20	100.0

<sup>1</sup>Intra-residue restraints, <sup>2</sup>Sequential restraints, <sup>3</sup>Medium range restraints, <sup>4</sup>Long range restraints,

<sup>5</sup>Inter-chain restraints, <sup>6</sup> Number of models with violations

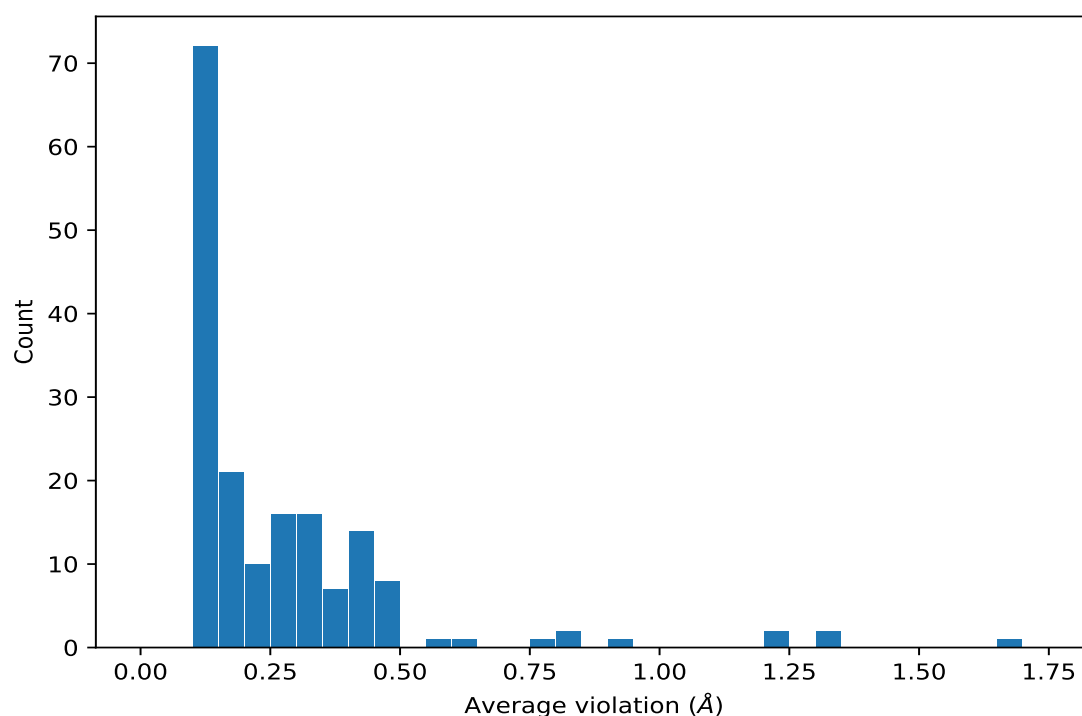
### 9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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

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

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



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

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

Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	20	1.68	0.28	1.66
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	20	1.32	0.15	1.33
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	20	1.31	0.23	1.31
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	20	1.22	0.35	1.07
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	20	1.22	0.93	0.95
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	20	0.91	0.23	0.93
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	20	0.84	0.08	0.84
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	20	0.84	0.24	0.84
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	20	0.75	0.11	0.74
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	20	0.62	0.36	0.63
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	20	0.6	0.12	0.6
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	20	0.49	0.16	0.52
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	20	0.49	0.18	0.44
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	20	0.46	0.17	0.52
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	20	0.39	0.21	0.3
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	20	0.39	0.09	0.38

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Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	20	0.36	0.06	0.36
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	20	0.36	0.06	0.36
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	20	0.3	0.07	0.29
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	20	0.27	0.09	0.24
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	20	0.25	0.08	0.25
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	20	0.18	0.06	0.17
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	20	0.18	0.04	0.18
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	19	0.12	0.01	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	19	0.12	0.01	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	19	0.12	0.01	0.12
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	18	0.26	0.06	0.26
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	18	0.26	0.06	0.26
(1,141)	1:8:A:GLY:H	1:37:A:CYS:H	18	0.26	0.06	0.26
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	18	0.22	0.04	0.24
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	18	0.22	0.04	0.24
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	18	0.22	0.04	0.24
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	18	0.22	0.04	0.24
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	17	0.3	0.1	0.28
(1,37)	1:4:A:ASN:HB2	1:19:A:TYR:H	17	0.3	0.1	0.28
(1,37)	1:4:A:ASN:HB3	1:19:A:TYR:H	17	0.3	0.1	0.28
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	16	0.18	0.05	0.16
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	16	0.15	0.04	0.14
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	16	0.15	0.04	0.14
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	16	0.15	0.04	0.14
(1,88)	1:28:A:LEU:HD21	1:55:A:VAL:H	16	0.15	0.04	0.14
(1,88)	1:28:A:LEU:HD22	1:55:A:VAL:H	16	0.15	0.04	0.14
(1,88)	1:28:A:LEU:HD23	1:55:A:VAL:H	16	0.15	0.04	0.14
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	15	0.32	0.1	0.29
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	15	0.32	0.1	0.29
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	15	0.32	0.1	0.29
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG11	15	0.32	0.1	0.29
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG12	15	0.32	0.1	0.29
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG13	15	0.32	0.1	0.29
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	12	0.14	0.03	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	12	0.14	0.03	0.13
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	12	0.13	0.02	0.12
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	11	0.46	0.09	0.43
(1,33)	1:71:A:LYS:HD2	1:72:A:VAL:H	11	0.46	0.09	0.43
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	11	0.11	0.01	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	11	0.11	0.01	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	11	0.11	0.01	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	11	0.11	0.01	0.12

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Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	11	0.11	0.01	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	11	0.11	0.01	0.12
(1,376)	1:106:A:LEU:HD11	1:123:A:TYR:HB2	10	0.17	0.06	0.15
(1,376)	1:106:A:LEU:HD12	1:123:A:TYR:HB2	10	0.17	0.06	0.15
(1,376)	1:106:A:LEU:HD13	1:123:A:TYR:HB2	10	0.17	0.06	0.15
(1,376)	1:28:A:LEU:HD11	1:54:A:TYR:HB3	10	0.17	0.06	0.15
(1,376)	1:28:A:LEU:HD12	1:54:A:TYR:HB3	10	0.17	0.06	0.15
(1,376)	1:28:A:LEU:HD13	1:54:A:TYR:HB3	10	0.17	0.06	0.15
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	9	0.11	0.01	0.1
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	9	0.11	0.01	0.1
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	9	0.11	0.01	0.1
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	8	0.19	0.04	0.18
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	8	0.15	0.02	0.14
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG21	7	0.33	0.1	0.36
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG22	7	0.33	0.1	0.36
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG23	7	0.33	0.1	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG21	7	0.33	0.1	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG22	7	0.33	0.1	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG23	7	0.33	0.1	0.36
(3,20)	1:72:A:VAL:N	1:68:A:ASP:O	7	0.16	0.04	0.17
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD2	7	0.13	0.01	0.14
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD3	7	0.13	0.01	0.14
(2,5)	1:79:A:GLN:H	1:56:A:GLY:O	7	0.13	0.01	0.13
(1,291)	1:103:A:TRP:HD1	1:120:A:VAL:HB	6	0.41	0.11	0.47
(1,291)	1:53:A:GLN:HB3	1:53:A:GLN:HE21	6	0.41	0.11	0.47
(1,167)	1:36:A:ARG:H	1:37:A:CYS:HB3	6	0.25	0.12	0.21
(1,167)	1:7:A:CYS:HB2	1:36:A:ARG:H	6	0.25	0.12	0.21
(1,167)	1:7:A:CYS:HB3	1:36:A:ARG:H	6	0.25	0.12	0.21
(1,982)	1:42:A:GLN:H	1:42:A:GLN:HE22	6	0.12	0.01	0.12
(1,839)	1:91:A:THR:HG21	1:92:A:LEU:H	6	0.11	0.01	0.11
(1,839)	1:91:A:THR:HG22	1:92:A:LEU:H	6	0.11	0.01	0.11
(1,839)	1:91:A:THR:HG23	1:92:A:LEU:H	6	0.11	0.01	0.11
(1,384)	1:6:A:VAL:HA	1:6:A:VAL:HB	5	0.37	0.12	0.43
(1,384)	1:116:A:GLU:HA	1:116:A:GLU:HB2	5	0.37	0.12	0.43
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG21	5	0.21	0.1	0.2
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG22	5	0.21	0.1	0.2
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG23	5	0.21	0.1	0.2
(1,367)	1:93:A:ARG:HB3	1:93:A:ARG:HD2	5	0.17	0.04	0.19
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD11	5	0.16	0.05	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD12	5	0.16	0.05	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD13	5	0.16	0.05	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD21	5	0.16	0.05	0.15

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Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD22	5	0.16	0.05	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD23	5	0.16	0.05	0.15
(1,496)	1:44:A:GLN:HA	1:44:A:GLN:HE22	5	0.12	0.01	0.13
(1,392)	1:110:A:GLU:HA	1:110:A:GLU:HG2	4	0.26	0.05	0.29
(1,392)	1:110:A:GLU:HA	1:110:A:GLU:HG3	4	0.26	0.05	0.29
(1,392)	1:53:A:GLN:HA	1:53:A:GLN:HG2	4	0.26	0.05	0.29
(1,392)	1:53:A:GLN:HA	1:53:A:GLN:HG3	4	0.26	0.05	0.29
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD21	4	0.14	0.05	0.12
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD22	4	0.14	0.05	0.12
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD23	4	0.14	0.05	0.12
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG2	4	0.12	0.01	0.12
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG3	4	0.12	0.01	0.12
(1,69)	1:19:A:TYR:HB2	1:22:A:GLY:H	4	0.11	0.02	0.11
(1,69)	1:20:A:ASP:HB2	1:22:A:GLY:H	4	0.11	0.02	0.11
(1,566)	1:93:A:ARG:HD3	1:94:A:GLU:H	4	0.11	0.0	0.11
(1,1556)	1:79:A:GLN:H	1:79:A:GLN:HB3	4	0.11	0.01	0.12
(1,86)	1:93:A:ARG:H	1:102:A:ILE:HG13	3	0.46	0.05	0.47
(1,86)	1:63:A:ILE:HG12	1:93:A:ARG:H	3	0.46	0.05	0.47
(1,86)	1:63:A:ILE:HG13	1:93:A:ARG:H	3	0.46	0.05	0.47
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG11	3	0.41	0.14	0.47
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG12	3	0.41	0.14	0.47
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG13	3	0.41	0.14	0.47
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG11	3	0.41	0.14	0.47
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG12	3	0.41	0.14	0.47
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG13	3	0.41	0.14	0.47
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG11	3	0.41	0.14	0.47
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG12	3	0.41	0.14	0.47
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG13	3	0.41	0.14	0.47
(1,476)	1:40:A:VAL:HG11	1:69:A:LYS:HG3	3	0.41	0.14	0.47
(1,476)	1:40:A:VAL:HG12	1:69:A:LYS:HG3	3	0.41	0.14	0.47
(1,476)	1:40:A:VAL:HG13	1:69:A:LYS:HG3	3	0.41	0.14	0.47
(1,17)	1:105:A:ALA:HB1	1:107:A:SER:H	3	0.33	0.14	0.42
(1,17)	1:105:A:ALA:HB2	1:107:A:SER:H	3	0.33	0.14	0.42
(1,17)	1:105:A:ALA:HB3	1:107:A:SER:H	3	0.33	0.14	0.42
(1,399)	1:93:A:ARG:HD2	1:102:A:ILE:HG12	3	0.21	0.07	0.22
(1,399)	1:84:A:GLU:HB3	1:88:A:TYR:HB3	3	0.21	0.07	0.22
(1,302)	1:40:A:VAL:HA	1:44:A:GLN:HE21	3	0.19	0.05	0.22
(1,302)	1:29:A:ILE:HA	1:30:A:PHE:HD1	3	0.19	0.05	0.22
(1,302)	1:29:A:ILE:HA	1:30:A:PHE:HD2	3	0.19	0.05	0.22
(2,4)	1:101:A:ALA:H	1:76:A:ALA:O	3	0.13	0.02	0.14
(1,1877)	1:13:A:GLN:HE21	1:14:A:ASP:HB2	3	0.13	0.02	0.12
(1,1877)	1:13:A:GLN:HE21	1:14:A:ASP:HB3	3	0.13	0.02	0.12

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Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(1,1877)	1:13:A:GLN:HE22	1:14:A:ASP:HB2	3	0.13	0.02	0.12
(1,1877)	1:13:A:GLN:HE22	1:14:A:ASP:HB3	3	0.13	0.02	0.12
(1,316)	1:39:A:ASN:HD22	1:42:A:GLN:HB2	2	0.37	0.01	0.37
(1,218)	1:79:A:GLN:HE21	1:103:A:TRP:HD1	2	0.32	0.02	0.32
(1,57)	1:14:A:ASP:HB3	1:17:A:ALA:H	2	0.21	0.01	0.21
(1,22)	1:35:A:PRO:HG2	1:36:A:ARG:H	2	0.14	0.02	0.14
(1,409)	1:69:A:LYS:HG2	1:69:A:LYS:HD2	2	0.12	0.01	0.12
(1,409)	1:69:A:LYS:HG2	1:69:A:LYS:HD3	2	0.12	0.01	0.12
(1,2000)	1:45:A:GLU:HB2	1:46:A:VAL:H	2	0.12	0.0	0.12
(1,2000)	1:45:A:GLU:HB3	1:46:A:VAL:H	2	0.12	0.0	0.12
(1,2106)	1:78:A:VAL:HG11	1:103:A:TRP:H	2	0.12	0.0	0.12
(1,2106)	1:78:A:VAL:HG12	1:103:A:TRP:H	2	0.12	0.0	0.12
(1,2106)	1:78:A:VAL:HG13	1:103:A:TRP:H	2	0.12	0.0	0.12
(1,2106)	1:78:A:VAL:HG21	1:103:A:TRP:H	2	0.12	0.0	0.12
(1,2106)	1:78:A:VAL:HG22	1:103:A:TRP:H	2	0.12	0.0	0.12
(1,2106)	1:78:A:VAL:HG23	1:103:A:TRP:H	2	0.12	0.0	0.12
(1,1298)	1:75:A:LEU:H	1:75:A:LEU:HG	2	0.12	0.02	0.12
(1,2149)	1:93:A:ARG:HB2	1:94:A:GLU:H	2	0.11	0.01	0.11
(1,2149)	1:93:A:ARG:HB3	1:94:A:GLU:H	2	0.11	0.01	0.11
(1,868)	1:102:A:ILE:HG21	1:121:A:ASP:H	2	0.11	0.0	0.11
(1,868)	1:102:A:ILE:HG22	1:121:A:ASP:H	2	0.11	0.0	0.11
(1,868)	1:102:A:ILE:HG23	1:121:A:ASP:H	2	0.11	0.0	0.11
(1,1353)	1:8:A:GLY:H	1:37:A:CYS:H	2	0.11	0.0	0.11
(1,1476)	1:29:A:ILE:HG21	1:56:A:GLY:HA2	2	0.11	0.0	0.11
(1,1476)	1:29:A:ILE:HG22	1:56:A:GLY:HA2	2	0.11	0.0	0.11
(1,1476)	1:29:A:ILE:HG23	1:56:A:GLY:HA2	2	0.11	0.0	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD11	2	0.11	0.0	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD12	2	0.11	0.0	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD13	2	0.11	0.0	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD21	2	0.11	0.0	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD22	2	0.11	0.0	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD23	2	0.11	0.0	0.11
(1,956)	1:12:A:GLY:HA3	1:13:A:GLN:HE21	2	0.1	0.0	0.1

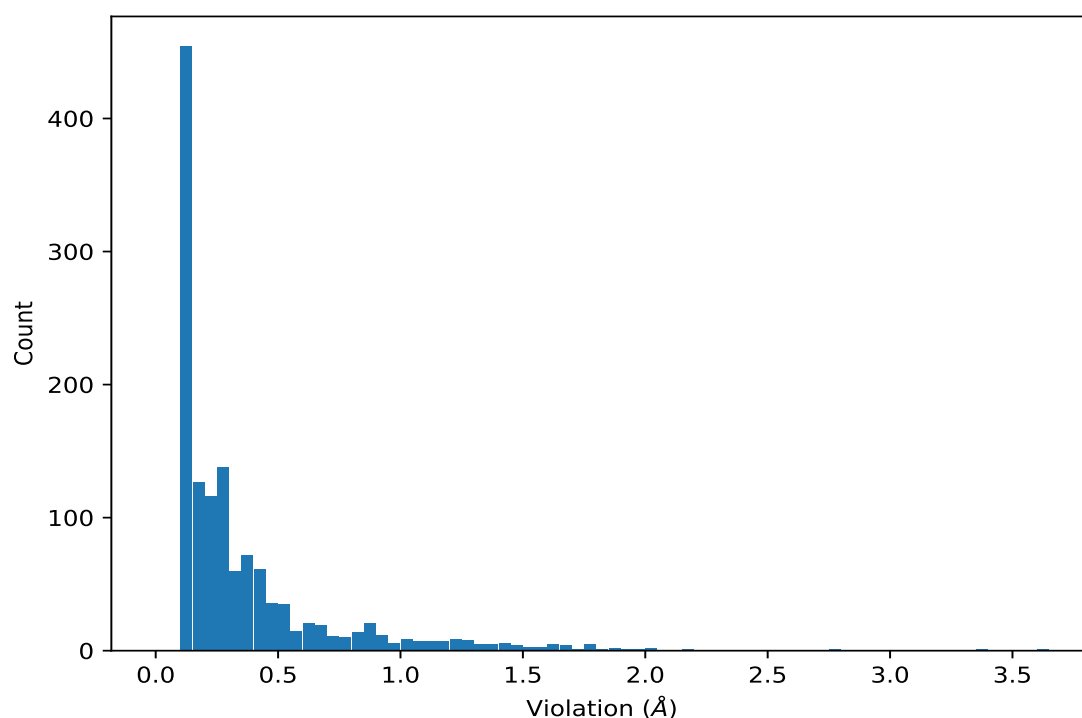
<sup>1</sup>Number of violated models, <sup>2</sup>Standard deviation

## 9.5 All violated distance restraints [i](#)

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

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





### 9.5.2 Table : All distance violations [i](#)

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

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	18	3.64
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	20	3.36
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	15	2.79
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	2	2.19
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	12	2.02
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	13	2.01
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	9	1.95
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	9	1.93
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	17	1.87
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	10	1.85
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	19	1.82
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	4	1.79
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	3	1.77
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	18	1.77
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	6	1.75
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	19	1.75

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	3	1.72
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	15	1.67
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	6	1.66
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	16	1.66
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	18	1.65
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	20	1.64
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	1	1.62
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	7	1.62
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	18	1.6
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	3	1.6
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	5	1.59
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	10	1.59
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	18	1.57
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	1	1.54
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	20	1.53
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	13	1.5
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	2	1.49
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	15	1.49
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	8	1.45
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	11	1.45
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	2	1.43
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	17	1.42
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	10	1.41
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	7	1.4
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	17	1.4
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	5	1.4
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	11	1.39
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	3	1.36
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	19	1.36
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	15	1.35
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	20	1.35
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	3	1.33
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	9	1.32
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	4	1.32
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	13	1.31
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	14	1.3
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	14	1.29
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	16	1.29
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	10	1.29
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	17	1.28
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	5	1.27
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	15	1.26

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	8	1.26
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	10	1.26
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	14	1.24
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	7	1.24
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	15	1.24
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	2	1.24
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	12	1.23
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	17	1.21
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	19	1.21
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	17	1.2
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	6	1.2
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	11	1.19
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	12	1.19
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	7	1.19
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	12	1.18
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	16	1.17
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	11	1.17
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	14	1.16
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	1	1.14
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	11	1.14
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	1	1.12
(3,10)	1:57:A:VAL:N	1:28:A:LEU:O	8	1.12
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	9	1.11
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	20	1.11
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	17	1.1
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	12	1.08
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	12	1.07
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	5	1.07
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	14	1.07
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	19	1.05
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	16	1.05
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	16	1.05
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	4	1.04
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	11	1.04
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	19	1.04
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	12	1.04
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	19	1.03
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	10	1.02
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	2	1.02
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	6	1.01
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	9	1.0
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	9	0.99

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	17	0.98
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	7	0.98
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	9	0.98
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	7	0.97
(3,25)	1:89:A:ILE:N	1:85:A:GLU:O	4	0.95
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	1	0.94
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	8	0.93
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	13	0.93
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	14	0.93
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	20	0.92
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	8	0.91
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	15	0.91
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	1	0.9
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	4	0.9
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	1	0.9
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	5	0.9
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	13	0.9
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	6	0.89
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	20	0.89
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	11	0.89
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	4	0.89
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	13	0.89
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	20	0.89
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	4	0.88
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	4	0.88
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	3	0.88
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	14	0.88
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	14	0.87
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	8	0.87
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	15	0.87
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	20	0.87
(3,24)	1:90:A:ASP:N	1:86:A:GLN:O	6	0.86
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	6	0.86
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	1	0.85
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	3	0.85
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	10	0.85
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	4	0.85
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	8	0.85
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	18	0.84
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	10	0.84
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	19	0.84
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	2	0.83

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	14	0.83
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	15	0.83
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	15	0.82
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	13	0.82
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	2	0.82
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	4	0.82
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	8	0.82
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	16	0.81
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	5	0.81
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	19	0.8
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	12	0.79
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	18	0.79
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	9	0.77
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	13	0.77
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	9	0.77
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	5	0.77
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	7	0.77
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	12	0.76
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	7	0.75
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	20	0.75
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	1	0.74
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	8	0.74
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	6	0.74
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	11	0.73
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	18	0.73
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	8	0.72
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	18	0.71
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	12	0.71
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	10	0.7
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	3	0.7
(1,427)	1:29:A:ILE:HA	1:57:A:VAL:HB	7	0.7
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	18	0.69
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	16	0.69
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	16	0.69
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	6	0.68
(3,21)	1:93:A:ARG:N	1:89:A:ILE:O	12	0.68
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	9	0.68
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	9	0.68
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	3	0.67
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	16	0.67
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	14	0.67
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	15	0.67

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	15	0.67
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	5	0.66
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	19	0.66
(3,8)	1:77:A:ALA:N	1:54:A:TYR:O	11	0.66
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	9	0.65
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	20	0.65
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	1	0.65
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	7	0.65
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	17	0.64
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	13	0.63
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	10	0.63
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	17	0.62
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	7	0.62
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	7	0.62
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	7	0.62
(1,33)	1:71:A:LYS:HD2	1:72:A:VAL:H	17	0.62
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	5	0.61
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	7	0.61
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	10	0.61
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	13	0.61
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	16	0.61
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	18	0.61
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	2	0.6
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	17	0.6
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	9	0.6
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	4	0.6
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	11	0.6
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	9	0.6
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	12	0.6
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	10	0.59
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	18	0.59
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	4	0.59
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	5	0.58
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	2	0.58
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	1	0.58
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	6	0.57
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	7	0.57
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	19	0.57
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	7	0.56
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	16	0.56
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	10	0.55
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	18	0.55

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	18	0.55
(1,33)	1:71:A:LYS:HD2	1:72:A:VAL:H	10	0.55
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	2	0.54
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	2	0.54
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	14	0.54
(3,12)	1:95:A:ALA:N	1:91:A:THR:O	17	0.54
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	11	0.54
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	2	0.54
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	6	0.54
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	15	0.54
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG11	8	0.54
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG12	8	0.54
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG13	8	0.54
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG11	8	0.54
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG12	8	0.54
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG13	8	0.54
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG11	8	0.54
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG12	8	0.54
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG13	8	0.54
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	13	0.53
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	5	0.53
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	14	0.52
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	19	0.52
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	14	0.52
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	7	0.51
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	14	0.51
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	20	0.51
(1,86)	1:63:A:ILE:HG12	1:93:A:ARG:H	11	0.51
(1,86)	1:63:A:ILE:HG13	1:93:A:ARG:H	11	0.51
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	1	0.5
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	8	0.5
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	13	0.5
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	10	0.5
(1,37)	1:4:A:ASN:HB2	1:19:A:TYR:H	2	0.5
(1,37)	1:4:A:ASN:HB3	1:19:A:TYR:H	2	0.5
(1,37)	1:4:A:ASN:HB2	1:19:A:TYR:H	9	0.5
(1,37)	1:4:A:ASN:HB3	1:19:A:TYR:H	9	0.5
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	13	0.49
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	11	0.49
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	1	0.49
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	3	0.49
(1,167)	1:7:A:CYS:HB2	1:36:A:ARG:H	16	0.49

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,167)	1:7:A:CYS:HB3	1:36:A:ARG:H	16	0.49
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	18	0.48
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	3	0.48
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	13	0.48
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	18	0.48
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	14	0.48
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	3	0.48
(1,291)	1:103:A:TRP:HD1	1:120:A:VAL:HB	14	0.48
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	3	0.47
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	3	0.47
(3,9)	1:55:A:VAL:N	1:26:A:GLY:O	8	0.47
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG11	18	0.47
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG12	18	0.47
(1,476)	1:38:A:VAL:HG11	1:40:A:VAL:HG13	18	0.47
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG11	18	0.47
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG12	18	0.47
(1,476)	1:38:A:VAL:HG12	1:40:A:VAL:HG13	18	0.47
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG11	18	0.47
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG12	18	0.47
(1,476)	1:38:A:VAL:HG13	1:40:A:VAL:HG13	18	0.47
(1,291)	1:53:A:GLN:HB3	1:53:A:GLN:HE21	9	0.47
(1,291)	1:103:A:TRP:HD1	1:120:A:VAL:HB	12	0.47
(1,291)	1:103:A:TRP:HD1	1:120:A:VAL:HB	13	0.47
(1,86)	1:93:A:ARG:H	1:102:A:ILE:HG13	16	0.47
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	20	0.46
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	5	0.46
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	14	0.46
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	5	0.46
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	2	0.45
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	18	0.45
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	9	0.45
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	1	0.44
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	1	0.44
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	1	0.44
(1,356)	1:57:A:VAL:HG21	1:77:A:ALA:HB1	1	0.44
(1,356)	1:57:A:VAL:HG21	1:77:A:ALA:HB2	1	0.44
(1,356)	1:57:A:VAL:HG21	1:77:A:ALA:HB3	1	0.44
(1,356)	1:57:A:VAL:HG22	1:77:A:ALA:HB1	1	0.44
(1,356)	1:57:A:VAL:HG22	1:77:A:ALA:HB2	1	0.44
(1,356)	1:57:A:VAL:HG22	1:77:A:ALA:HB3	1	0.44
(1,356)	1:57:A:VAL:HG23	1:77:A:ALA:HB1	1	0.44
(1,356)	1:57:A:VAL:HG23	1:77:A:ALA:HB2	1	0.44

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,356)	1:57:A:VAL:HG23	1:77:A:ALA:HB3	1	0.44
(1,17)	1:105:A:ALA:HB1	1:107:A:SER:H	3	0.44
(1,17)	1:105:A:ALA:HB2	1:107:A:SER:H	3	0.44
(1,17)	1:105:A:ALA:HB3	1:107:A:SER:H	3	0.44
(3,28)	1:115:A:ARG:N	1:113:A:PRO:O	5	0.43
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	9	0.43
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	16	0.43
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	7	0.43
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	9	0.43
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	2	0.43
(1,384)	1:6:A:VAL:HA	1:6:A:VAL:HB	3	0.43
(1,384)	1:116:A:GLU:HA	1:116:A:GLU:HB2	11	0.43
(1,384)	1:6:A:VAL:HA	1:6:A:VAL:HB	20	0.43
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	3	0.43
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	3	0.42
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	4	0.42
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	15	0.42
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	2	0.42
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	4	0.42
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	11	0.42
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	19	0.42
(1,17)	1:105:A:ALA:HB1	1:107:A:SER:H	12	0.42
(1,17)	1:105:A:ALA:HB2	1:107:A:SER:H	12	0.42
(1,17)	1:105:A:ALA:HB3	1:107:A:SER:H	12	0.42
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	8	0.41
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	12	0.41
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	1	0.41
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	11	0.41
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	13	0.41
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	4	0.41
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	12	0.41
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	1	0.41
(1,384)	1:116:A:GLU:HA	1:116:A:GLU:HB2	6	0.41
(1,291)	1:103:A:TRP:HD1	1:120:A:VAL:HB	7	0.41
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	16	0.41
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	20	0.4
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	6	0.4
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	15	0.4
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	19	0.4
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	6	0.4
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	15	0.4
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG21	2	0.4

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG22	2	0.4
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG23	2	0.4
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG21	2	0.4
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG22	2	0.4
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG23	2	0.4
(1,86)	1:63:A:ILE:HG12	1:93:A:ARG:H	6	0.4
(1,86)	1:63:A:ILE:HG13	1:93:A:ARG:H	6	0.4
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	13	0.4
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	16	0.39
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	1	0.39
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	5	0.39
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	15	0.39
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	3	0.39
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	4	0.39
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	20	0.39
(3,13)	1:66:A:VAL:N	1:62:A:ASP:O	3	0.39
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG21	10	0.39
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG22	10	0.39
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG23	10	0.39
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG21	10	0.39
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG22	10	0.39
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG23	10	0.39
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	17	0.38
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	4	0.38
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	17	0.38
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	8	0.38
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	12	0.38
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG21	14	0.38
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG22	14	0.38
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG23	14	0.38
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	20	0.38
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	20	0.38
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	20	0.38
(1,316)	1:39:A:ASN:HD22	1:42:A:GLN:HB2	7	0.38
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG21	12	0.38
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG22	12	0.38
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG23	12	0.38
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG21	12	0.38
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG22	12	0.38
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG23	12	0.38
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	11	0.37
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	20	0.37

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	15	0.37
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	16	0.37
(1,37)	1:4:A:ASN:HB2	1:19:A:TYR:H	19	0.37
(1,37)	1:4:A:ASN:HB3	1:19:A:TYR:H	19	0.37
(3,29)	1:84:A:GLU:N	1:81:A:HIS:O	7	0.36
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	10	0.36
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	11	0.36
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	13	0.36
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	10	0.36
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	11	0.36
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	13	0.36
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	9	0.36
(1,316)	1:39:A:ASN:HD22	1:42:A:GLN:HB2	15	0.36
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG21	11	0.36
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG22	11	0.36
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG23	11	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG21	11	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG22	11	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG23	11	0.36
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG21	19	0.36
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG22	19	0.36
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG23	19	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG21	19	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG22	19	0.36
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG23	19	0.36
(1,141)	1:8:A:GLY:H	1:37:A:CYS:H	20	0.36
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	7	0.36
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	2	0.35
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	14	0.35
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	2	0.35
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	14	0.35
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	12	0.35
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	19	0.35
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	2	0.35
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	14	0.35
(1,162)	1:6:A:VAL:H	1:18:A:ALA:HB1	11	0.35
(1,162)	1:6:A:VAL:H	1:18:A:ALA:HB2	11	0.35
(1,162)	1:6:A:VAL:H	1:18:A:ALA:HB3	11	0.35
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	12	0.34
(1,218)	1:79:A:GLN:HE21	1:103:A:TRP:HD1	20	0.34
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	1	0.33
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	4	0.33

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	4	0.33
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	7	0.33
(1,141)	1:8:A:GLY:H	1:37:A:CYS:H	7	0.33
(1,141)	1:8:A:GLY:H	1:37:A:CYS:H	8	0.33
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	14	0.33
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	14	0.33
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	12	0.33
(1,33)	1:71:A:LYS:HD3	1:72:A:VAL:H	6	0.33
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	7	0.32
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	20	0.32
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	8	0.32
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	9	0.32
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	3	0.32
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	19	0.32
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	7	0.32
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	20	0.32
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	9	0.32
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	6	0.32
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	6	0.32
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	6	0.32
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	13	0.32
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	13	0.32
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	1	0.31
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	16	0.31
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	19	0.31
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	20	0.31
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	13	0.31
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	1	0.31
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	16	0.31
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	4	0.31
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	15	0.31
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	16	0.31
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG21	3	0.31
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG22	3	0.31
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG23	3	0.31
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG21	3	0.31
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG22	3	0.31
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG23	3	0.31
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	9	0.3
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	6	0.3
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	6	0.3
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	8	0.3

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	3	0.3
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	3	0.3
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	3	0.3
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG11	13	0.3
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG12	13	0.3
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG13	13	0.3
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG11	15	0.3
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG12	15	0.3
(1,411)	1:42:A:GLN:HB2	1:46:A:VAL:HG13	15	0.3
(1,392)	1:110:A:GLU:HA	1:110:A:GLU:HG2	7	0.3
(1,392)	1:110:A:GLU:HA	1:110:A:GLU:HG3	7	0.3
(1,218)	1:79:A:GLN:HE21	1:103:A:TRP:HD1	11	0.3
(1,141)	1:8:A:GLY:H	1:37:A:CYS:H	6	0.3
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	10	0.3
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	10	0.29
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	16	0.29
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	5	0.29
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	11	0.29
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	20	0.29
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	5	0.29
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	8	0.29
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	8	0.29
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	8	0.29
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	10	0.29
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	10	0.29
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	10	0.29
(1,399)	1:84:A:GLU:HB3	1:88:A:TYR:HB3	5	0.29
(1,392)	1:110:A:GLU:HA	1:110:A:GLU:HG2	9	0.29
(1,392)	1:110:A:GLU:HA	1:110:A:GLU:HG3	9	0.29
(1,190)	1:52:A:LEU:HD11	1:53:A:GLN:H	1	0.29
(1,190)	1:52:A:LEU:HD12	1:53:A:GLN:H	1	0.29
(1,190)	1:52:A:LEU:HD13	1:53:A:GLN:H	1	0.29
(1,190)	1:52:A:LEU:HD21	1:53:A:GLN:H	1	0.29
(1,190)	1:52:A:LEU:HD22	1:53:A:GLN:H	1	0.29
(1,190)	1:52:A:LEU:HD23	1:53:A:GLN:H	1	0.29
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	13	0.28
(3,22)	1:91:A:THR:N	1:87:A:LEU:O	8	0.28
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	2	0.28
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	12	0.28
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	11	0.28
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	17	0.28
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	6	0.28

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	17	0.28
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	10	0.28
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	17	0.28
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	14	0.28
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	14	0.28
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	14	0.28
(1,392)	1:53:A:GLN:HA	1:53:A:GLN:HG2	12	0.28
(1,392)	1:53:A:GLN:HA	1:53:A:GLN:HG3	12	0.28
(1,167)	1:36:A:ARG:H	1:37:A:CYS:HB3	20	0.28
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	4	0.28
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	4	0.28
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	5	0.28
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	8	0.28
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	17	0.28
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	4	0.27
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	8	0.27
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	19	0.27
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	7	0.27
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	14	0.27
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	11	0.27
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	18	0.27
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	8	0.27
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	19	0.27
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	8	0.27
(3,6)	1:56:A:GLY:N	1:77:A:ALA:O	13	0.27
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	16	0.27
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	16	0.27
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	16	0.27
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	14	0.27
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	14	0.27
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	14	0.27
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	14	0.27
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	17	0.27
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	17	0.27
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	17	0.27
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	17	0.27
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	5	0.27
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	5	0.27
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	15	0.27
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	15	0.27
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	1	0.27
(1,37)	1:4:A:ASN:HB2	1:19:A:TYR:H	16	0.27

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,37)	1:4:A:ASN:HB3	1:19:A:TYR:H	16	0.27
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	6	0.26
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	2	0.26
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	5	0.26
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	6	0.26
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	12	0.26
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	6	0.26
(1,376)	1:106:A:LEU:HD11	1:123:A:TYR:HB2	2	0.26
(1,376)	1:106:A:LEU:HD12	1:123:A:TYR:HB2	2	0.26
(1,376)	1:106:A:LEU:HD13	1:123:A:TYR:HB2	2	0.26
(1,376)	1:106:A:LEU:HD11	1:123:A:TYR:HB2	15	0.26
(1,376)	1:106:A:LEU:HD12	1:123:A:TYR:HB2	15	0.26
(1,376)	1:106:A:LEU:HD13	1:123:A:TYR:HB2	15	0.26
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD11	15	0.26
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD12	15	0.26
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD13	15	0.26
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD21	15	0.26
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD22	15	0.26
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD23	15	0.26
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	10	0.26
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	10	0.26
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	10	0.26
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	10	0.26
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	17	0.26
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	17	0.26
(1,16)	1:42:A:GLN:HA	1:44:A:GLN:H	8	0.26
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	15	0.25
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	8	0.25
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	14	0.25
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	5	0.25
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	8	0.25
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	4	0.25
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	17	0.25
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	14	0.25
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	15	0.25
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	12	0.25
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	14	0.25
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	16	0.25
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	4	0.25
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	4	0.25
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	4	0.25
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	12	0.25

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	12	0.25
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	12	0.25
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	19	0.25
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	19	0.25
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	19	0.25
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	5	0.25
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	5	0.25
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	5	0.25
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	5	0.25
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	7	0.25
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	7	0.25
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	7	0.25
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	7	0.25
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	13	0.25
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	13	0.25
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	13	0.25
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	13	0.25
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	16	0.25
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	16	0.25
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	16	0.25
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	16	0.25
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	1	0.25
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	1	0.25
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	3	0.25
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	13	0.25
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	20	0.25
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	11	0.24
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	12	0.24
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	3	0.24
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	18	0.24
(3,16)	1:69:A:LYS:N	1:65:A:ASP:O	6	0.24
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	11	0.24
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	1	0.24
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	16	0.24
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	3	0.24
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG11	17	0.24
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG12	17	0.24
(1,411)	1:45:A:GLU:HB2	1:46:A:VAL:HG13	17	0.24
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	1	0.24
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	1	0.24
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	1	0.24
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	1	0.24

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	4	0.24
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	4	0.24
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	4	0.24
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	4	0.24
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	6	0.24
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	6	0.24
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	6	0.24
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	6	0.24
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	10	0.24
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	10	0.24
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	9	0.24
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	9	0.24
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	9	0.24
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	7	0.24
(3,20)	1:72:A:VAL:N	1:68:A:ASP:O	16	0.23
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	2	0.23
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	7	0.23
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	20	0.23
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG21	3	0.23
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG22	3	0.23
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG23	3	0.23
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD21	15	0.23
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD22	15	0.23
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD23	15	0.23
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	2	0.23
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	2	0.23
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	2	0.23
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	2	0.23
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	3	0.23
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	3	0.23
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	3	0.23
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	3	0.23
(1,302)	1:40:A:VAL:HA	1:44:A:GLN:HE21	16	0.23
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	5	0.22
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	16	0.22
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	11	0.22
(3,23)	1:92:A:LEU:N	1:88:A:TYR:O	5	0.22
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	17	0.22
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	20	0.22
(3,11)	1:92:A:LEU:N	1:88:A:TYR:O	5	0.22
(1,1338)	1:39:A:ASN:HB2	1:39:A:ASN:HD21	8	0.22
(1,453)	1:44:A:GLN:HG2	1:72:A:VAL:HG11	16	0.22

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,453)	1:44:A:GLN:HG2	1:72:A:VAL:HG12	16	0.22
(1,453)	1:44:A:GLN:HG2	1:72:A:VAL:HG13	16	0.22
(1,399)	1:84:A:GLU:HB3	1:88:A:TYR:HB3	7	0.22
(1,302)	1:29:A:ILE:HA	1:30:A:PHE:HD1	10	0.22
(1,302)	1:29:A:ILE:HA	1:30:A:PHE:HD2	10	0.22
(1,167)	1:36:A:ARG:H	1:37:A:CYS:HB3	6	0.22
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	2	0.22
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	2	0.22
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	19	0.22
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	19	0.22
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	19	0.22
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	19	0.22
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	19	0.22
(1,57)	1:14:A:ASP:HB3	1:17:A:ALA:H	13	0.22
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	6	0.21
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	18	0.21
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	17	0.21
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	5	0.21
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	18	0.21
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	6	0.21
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	13	0.21
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	8	0.21
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	8	0.21
(1,476)	1:40:A:VAL:HG11	1:69:A:LYS:HG3	5	0.21
(1,476)	1:40:A:VAL:HG12	1:69:A:LYS:HG3	5	0.21
(1,476)	1:40:A:VAL:HG13	1:69:A:LYS:HG3	5	0.21
(1,376)	1:106:A:LEU:HD11	1:123:A:TYR:HB2	19	0.21
(1,376)	1:106:A:LEU:HD12	1:123:A:TYR:HB2	19	0.21
(1,376)	1:106:A:LEU:HD13	1:123:A:TYR:HB2	19	0.21
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	18	0.21
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	18	0.21
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	15	0.21
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	20	0.21
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	3	0.2
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	1	0.2
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	11	0.2
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	5	0.2
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	5	0.2
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG21	20	0.2
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG22	20	0.2
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG23	20	0.2
(1,367)	1:93:A:ARG:HB3	1:93:A:ARG:HD2	17	0.2

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,367)	1:93:A:ARG:HB3	1:93:A:ARG:HD2	20	0.2
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	11	0.2
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	11	0.2
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	11	0.2
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	11	0.2
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	18	0.2
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	18	0.2
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	18	0.2
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	18	0.2
(1,167)	1:36:A:ARG:H	1:37:A:CYS:HB3	5	0.2
(1,116)	1:42:A:GLN:H	1:42:A:GLN:HG3	8	0.2
(1,57)	1:14:A:ASP:HB3	1:17:A:ALA:H	10	0.2
(1,37)	1:4:A:ASN:HB2	1:19:A:TYR:H	6	0.2
(1,37)	1:4:A:ASN:HB3	1:19:A:TYR:H	6	0.2
(1,20)	1:25:A:TYR:HB2	1:53:A:GLN:H	7	0.2
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	15	0.2
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	9	0.19
(3,20)	1:72:A:VAL:N	1:68:A:ASP:O	18	0.19
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	4	0.19
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	13	0.19
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	9	0.19
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	14	0.19
(1,491)	1:80:A:LEU:H	1:82:A:GLY:H	14	0.19
(1,367)	1:93:A:ARG:HB3	1:93:A:ARG:HD2	9	0.19
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	12	0.19
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	12	0.19
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	12	0.19
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	12	0.19
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	7	0.18
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	19	0.18
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	5	0.18
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	18	0.18
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	10	0.18
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	12	0.18
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	16	0.18
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	18	0.18
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	2	0.18
(3,3)	1:80:A:LEU:N	1:103:A:TRP:O	17	0.18
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	3	0.18
(1,477)	1:44:A:GLN:HG2	1:72:A:VAL:HG11	16	0.18
(1,477)	1:44:A:GLN:HG2	1:72:A:VAL:HG12	16	0.18
(1,477)	1:44:A:GLN:HG2	1:72:A:VAL:HG13	16	0.18

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,451)	1:28:A:LEU:HA	1:28:A:LEU:HD21	11	0.18
(1,451)	1:28:A:LEU:HA	1:28:A:LEU:HD22	11	0.18
(1,451)	1:28:A:LEU:HA	1:28:A:LEU:HD23	11	0.18
(1,392)	1:53:A:GLN:HA	1:53:A:GLN:HG2	14	0.18
(1,392)	1:53:A:GLN:HA	1:53:A:GLN:HG3	14	0.18
(1,376)	1:106:A:LEU:HD11	1:123:A:TYR:HB2	5	0.18
(1,376)	1:106:A:LEU:HD12	1:123:A:TYR:HB2	5	0.18
(1,376)	1:106:A:LEU:HD13	1:123:A:TYR:HB2	5	0.18
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	8	0.18
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	8	0.18
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	8	0.18
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	8	0.18
(1,249)	1:89:A:ILE:HB	1:117:A:PHE:HD1	1	0.18
(1,249)	1:89:A:ILE:HB	1:117:A:PHE:HD2	1	0.18
(1,167)	1:36:A:ARG:H	1:37:A:CYS:HB3	13	0.18
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	3	0.18
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	3	0.18
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	17	0.18
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	12	0.17
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	15	0.17
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	16	0.17
(3,20)	1:72:A:VAL:N	1:68:A:ASP:O	2	0.17
(3,20)	1:72:A:VAL:N	1:68:A:ASP:O	12	0.17
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	1	0.17
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	14	0.17
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	15	0.17
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	6	0.17
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	3	0.17
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	8	0.17
(1,372)	1:41:A:GLU:HA	1:41:A:GLU:HG3	8	0.17
(1,367)	1:93:A:ARG:HB3	1:93:A:ARG:HD2	13	0.17
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	15	0.17
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	15	0.17
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	15	0.17
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	15	0.17
(1,291)	1:53:A:GLN:HB3	1:53:A:GLN:HE21	20	0.17
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	3	0.17
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	3	0.17
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	3	0.17
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	3	0.17
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	7	0.17
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	12	0.17

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	6	0.16
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	11	0.16
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	4	0.16
(1,1877)	1:13:A:GLN:HE21	1:14:A:ASP:HB2	19	0.16
(1,1877)	1:13:A:GLN:HE21	1:14:A:ASP:HB3	19	0.16
(1,1877)	1:13:A:GLN:HE22	1:14:A:ASP:HB2	19	0.16
(1,1877)	1:13:A:GLN:HE22	1:14:A:ASP:HB3	19	0.16
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	15	0.16
(1,390)	1:67:A:VAL:HG21	1:96:A:LEU:HA	12	0.16
(1,390)	1:67:A:VAL:HG22	1:96:A:LEU:HA	12	0.16
(1,390)	1:67:A:VAL:HG23	1:96:A:LEU:HA	12	0.16
(1,376)	1:106:A:LEU:HD11	1:123:A:TYR:HB2	3	0.16
(1,376)	1:106:A:LEU:HD12	1:123:A:TYR:HB2	3	0.16
(1,376)	1:106:A:LEU:HD13	1:123:A:TYR:HB2	3	0.16
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	12	0.16
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	12	0.16
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	12	0.16
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	18	0.16
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	18	0.16
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	18	0.16
(1,22)	1:35:A:PRO:HG2	1:36:A:ARG:H	15	0.16
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	10	0.15
(3,27)	1:100:A:VAL:N	1:97:A:PRO:O	17	0.15
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	20	0.15
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	20	0.15
(3,18)	1:73:A:LEU:N	1:69:A:LYS:O	10	0.15
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	9	0.15
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	10	0.15
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	12	0.15
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	16	0.15
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	10	0.15
(3,14)	1:68:A:ASP:N	1:64:A:ALA:O	15	0.15
(2,5)	1:79:A:GLN:H	1:56:A:GLY:O	2	0.15
(2,4)	1:101:A:ALA:H	1:76:A:ALA:O	17	0.15
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD2	2	0.15
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD3	2	0.15
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD2	10	0.15
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD3	10	0.15
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	2	0.15
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	2	0.15
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	20	0.15
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	20	0.15

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD11	1	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD12	1	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD13	1	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD21	1	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD22	1	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD23	1	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD11	7	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD12	7	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD13	7	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD21	7	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD22	7	0.15
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD23	7	0.15
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	12	0.15
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	12	0.15
(1,88)	1:28:A:LEU:HD21	1:55:A:VAL:H	6	0.15
(1,88)	1:28:A:LEU:HD22	1:55:A:VAL:H	6	0.15
(1,88)	1:28:A:LEU:HD23	1:55:A:VAL:H	6	0.15
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	1	0.14
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	13	0.14
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	17	0.14
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	19	0.14
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	1	0.14
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	8	0.14
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	10	0.14
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	19	0.14
(2,5)	1:79:A:GLN:H	1:56:A:GLY:O	12	0.14
(2,4)	1:101:A:ALA:H	1:76:A:ALA:O	11	0.14
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD2	3	0.14
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD3	3	0.14
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD2	18	0.14
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD3	18	0.14
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	15	0.14
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	15	0.14
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	8	0.14
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	14	0.14
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	20	0.14
(1,982)	1:42:A:GLN:H	1:42:A:GLN:HE22	20	0.14
(1,962)	1:39:A:ASN:HD21	1:41:A:GLU:HA	8	0.14
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	3	0.14
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	3	0.14
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	3	0.14
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	4	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	4	0.14
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	4	0.14
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG2	5	0.14
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG3	5	0.14
(1,376)	1:28:A:LEU:HD11	1:54:A:TYR:HB3	20	0.14
(1,376)	1:28:A:LEU:HD12	1:54:A:TYR:HB3	20	0.14
(1,376)	1:28:A:LEU:HD13	1:54:A:TYR:HB3	20	0.14
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD11	20	0.14
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD12	20	0.14
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD13	20	0.14
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD21	20	0.14
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD22	20	0.14
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD23	20	0.14
(1,141)	1:123:A:TYR:HD1	1:124:A:VAL:H	16	0.14
(1,141)	1:123:A:TYR:HD2	1:124:A:VAL:H	16	0.14
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	4	0.14
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	4	0.14
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	4	0.14
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	13	0.14
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	13	0.14
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	13	0.14
(1,69)	1:20:A:ASP:HB2	1:22:A:GLY:H	10	0.14
(1,64)	1:39:A:ASN:HD21	1:41:A:GLU:HA	8	0.14
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	2	0.13
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	3	0.13
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	10	0.13
(3,20)	1:72:A:VAL:N	1:68:A:ASP:O	6	0.13
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	9	0.13
(3,19)	1:70:A:ALA:N	1:66:A:VAL:O	13	0.13
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	1	0.13
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	19	0.13
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	19	0.13
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	6	0.13
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	17	0.13
(2,5)	1:79:A:GLN:H	1:56:A:GLY:O	1	0.13
(2,5)	1:79:A:GLN:H	1:56:A:GLY:O	3	0.13
(2,5)	1:79:A:GLN:H	1:56:A:GLY:O	10	0.13
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	12	0.13
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	12	0.13
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	12	0.13
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	12	0.13
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	12	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	12	0.13
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	15	0.13
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	15	0.13
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	15	0.13
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	15	0.13
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	15	0.13
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	15	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	4	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	4	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	11	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	11	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	12	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	12	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	19	0.13
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	19	0.13
(1,1298)	1:75:A:LEU:H	1:75:A:LEU:HG	17	0.13
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	1	0.13
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	4	0.13
(1,982)	1:42:A:GLN:H	1:42:A:GLN:HE22	19	0.13
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	8	0.13
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	8	0.13
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	8	0.13
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	14	0.13
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	14	0.13
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	14	0.13
(1,496)	1:44:A:GLN:HA	1:44:A:GLN:HE22	1	0.13
(1,496)	1:44:A:GLN:HA	1:44:A:GLN:HE22	7	0.13
(1,496)	1:44:A:GLN:HA	1:44:A:GLN:HE22	20	0.13
(1,409)	1:69:A:LYS:HG2	1:69:A:LYS:HD2	5	0.13
(1,409)	1:69:A:LYS:HG2	1:69:A:LYS:HD3	5	0.13
(1,384)	1:116:A:GLU:HA	1:116:A:GLU:HB2	4	0.13
(1,376)	1:28:A:LEU:HD11	1:54:A:TYR:HB3	8	0.13
(1,376)	1:28:A:LEU:HD12	1:54:A:TYR:HB3	8	0.13
(1,376)	1:28:A:LEU:HD13	1:54:A:TYR:HB3	8	0.13
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE2	20	0.13
(1,312)	1:30:A:PHE:HD1	1:69:A:LYS:HE3	20	0.13
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE2	20	0.13
(1,312)	1:30:A:PHE:HD2	1:69:A:LYS:HE3	20	0.13
(1,194)	1:100:A:VAL:H	1:100:A:VAL:HG21	1	0.13
(1,194)	1:100:A:VAL:H	1:100:A:VAL:HG22	1	0.13
(1,194)	1:100:A:VAL:H	1:100:A:VAL:HG23	1	0.13
(1,167)	1:36:A:ARG:H	1:37:A:CYS:HB3	14	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	2	0.13
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	2	0.13
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	2	0.13
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	17	0.13
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	17	0.13
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	17	0.13
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	20	0.13
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	20	0.13
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	20	0.13
(1,37)	1:19:A:TYR:H	1:52:A:LEU:HG	14	0.13
(1,17)	1:105:A:ALA:HB1	1:107:A:SER:H	8	0.13
(1,17)	1:105:A:ALA:HB2	1:107:A:SER:H	8	0.13
(1,17)	1:105:A:ALA:HB3	1:107:A:SER:H	8	0.13
(3,20)	1:72:A:VAL:N	1:68:A:ASP:O	3	0.12
(3,20)	1:72:A:VAL:N	1:68:A:ASP:O	15	0.12
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	2	0.12
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	7	0.12
(3,15)	1:67:A:VAL:N	1:63:A:ILE:O	17	0.12
(3,7)	1:58:A:PHE:N	1:79:A:GLN:O	19	0.12
(2,5)	1:79:A:GLN:H	1:56:A:GLY:O	4	0.12
(1,2149)	1:93:A:ARG:HB2	1:94:A:GLU:H	5	0.12
(1,2149)	1:93:A:ARG:HB3	1:94:A:GLU:H	5	0.12
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD2	6	0.12
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD3	6	0.12
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD2	12	0.12
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD3	12	0.12
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD2	16	0.12
(1,2143)	1:92:A:LEU:H	1:93:A:ARG:HD3	16	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	10	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	10	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	10	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	10	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	10	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	10	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	11	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	11	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	11	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	11	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	11	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	11	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	19	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	19	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	19	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	19	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	19	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	19	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	20	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	20	0.12
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	20	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	20	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	20	0.12
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	20	0.12
(1,2106)	1:78:A:VAL:HG11	1:103:A:TRP:H	14	0.12
(1,2106)	1:78:A:VAL:HG12	1:103:A:TRP:H	14	0.12
(1,2106)	1:78:A:VAL:HG13	1:103:A:TRP:H	14	0.12
(1,2106)	1:78:A:VAL:HG21	1:103:A:TRP:H	14	0.12
(1,2106)	1:78:A:VAL:HG22	1:103:A:TRP:H	14	0.12
(1,2106)	1:78:A:VAL:HG23	1:103:A:TRP:H	14	0.12
(1,2106)	1:78:A:VAL:HG11	1:103:A:TRP:H	20	0.12
(1,2106)	1:78:A:VAL:HG12	1:103:A:TRP:H	20	0.12
(1,2106)	1:78:A:VAL:HG13	1:103:A:TRP:H	20	0.12
(1,2106)	1:78:A:VAL:HG21	1:103:A:TRP:H	20	0.12
(1,2106)	1:78:A:VAL:HG22	1:103:A:TRP:H	20	0.12
(1,2106)	1:78:A:VAL:HG23	1:103:A:TRP:H	20	0.12
(1,2000)	1:45:A:GLU:HB2	1:46:A:VAL:H	13	0.12
(1,2000)	1:45:A:GLU:HB3	1:46:A:VAL:H	13	0.12
(1,2000)	1:45:A:GLU:HB2	1:46:A:VAL:H	15	0.12
(1,2000)	1:45:A:GLU:HB3	1:46:A:VAL:H	15	0.12
(1,1877)	1:13:A:GLN:HE21	1:14:A:ASP:HB2	5	0.12
(1,1877)	1:13:A:GLN:HE21	1:14:A:ASP:HB3	5	0.12
(1,1877)	1:13:A:GLN:HE22	1:14:A:ASP:HB2	5	0.12
(1,1877)	1:13:A:GLN:HE22	1:14:A:ASP:HB3	5	0.12
(1,1556)	1:79:A:GLN:H	1:79:A:GLN:HB3	15	0.12
(1,1556)	1:79:A:GLN:H	1:79:A:GLN:HB3	20	0.12
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	10	0.12
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	10	0.12
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	17	0.12
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	17	0.12
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	2	0.12
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	2	0.12
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	2	0.12
(1,1194)	1:41:A:GLU:H	1:44:A:GLN:HE22	11	0.12
(1,1064)	1:34:A:SER:HA	1:37:A:CYS:H	5	0.12
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	5	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	6	0.12
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	13	0.12
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	18	0.12
(1,982)	1:42:A:GLN:H	1:42:A:GLN:HE22	4	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	5	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	5	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	5	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	6	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	6	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	6	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	13	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	13	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	13	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	16	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	16	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	16	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	19	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	19	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	19	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	20	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	20	0.12
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	20	0.12
(1,839)	1:91:A:THR:HG21	1:92:A:LEU:H	10	0.12
(1,839)	1:91:A:THR:HG22	1:92:A:LEU:H	10	0.12
(1,839)	1:91:A:THR:HG23	1:92:A:LEU:H	10	0.12
(1,839)	1:91:A:THR:HG21	1:92:A:LEU:H	18	0.12
(1,839)	1:91:A:THR:HG22	1:92:A:LEU:H	18	0.12
(1,839)	1:91:A:THR:HG23	1:92:A:LEU:H	18	0.12
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG2	20	0.12
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG3	20	0.12
(1,566)	1:93:A:ARG:HD3	1:94:A:GLU:H	16	0.12
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG21	1	0.12
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG22	1	0.12
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG23	1	0.12
(1,399)	1:93:A:ARG:HD2	1:102:A:ILE:HG12	13	0.12
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD21	7	0.12
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD22	7	0.12
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD23	7	0.12
(1,376)	1:28:A:LEU:HD11	1:54:A:TYR:HB3	4	0.12
(1,376)	1:28:A:LEU:HD12	1:54:A:TYR:HB3	4	0.12
(1,376)	1:28:A:LEU:HD13	1:54:A:TYR:HB3	4	0.12
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	1	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	1	0.12
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	1	0.12
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	8	0.12
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	8	0.12
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	8	0.12
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	10	0.12
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	10	0.12
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	10	0.12
(1,88)	1:28:A:LEU:HD21	1:55:A:VAL:H	16	0.12
(1,88)	1:28:A:LEU:HD22	1:55:A:VAL:H	16	0.12
(1,88)	1:28:A:LEU:HD23	1:55:A:VAL:H	16	0.12
(1,22)	1:35:A:PRO:HG2	1:36:A:ARG:H	9	0.12
(1,13)	1:73:A:LEU:HB2	1:74:A:SER:H	8	0.12
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	7	0.11
(3,26)	1:96:A:LEU:N	1:92:A:LEU:O	19	0.11
(3,17)	1:71:A:LYS:N	1:67:A:VAL:O	8	0.11
(3,2)	1:78:A:VAL:N	1:101:A:ALA:O	15	0.11
(2,4)	1:101:A:ALA:H	1:76:A:ALA:O	16	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD11	18	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD12	18	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD13	18	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD21	18	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD22	18	0.11
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD23	18	0.11
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	6	0.11
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	6	0.11
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	6	0.11
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	6	0.11
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	6	0.11
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	6	0.11
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	14	0.11
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	14	0.11
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	14	0.11
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	14	0.11
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	14	0.11
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	14	0.11
(1,1556)	1:79:A:GLN:H	1:79:A:GLN:HB3	11	0.11
(1,1476)	1:29:A:ILE:HG21	1:56:A:GLY:HA2	16	0.11
(1,1476)	1:29:A:ILE:HG22	1:56:A:GLY:HA2	16	0.11
(1,1476)	1:29:A:ILE:HG23	1:56:A:GLY:HA2	16	0.11
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE1	14	0.11
(1,1362)	1:25:A:TYR:HA	1:25:A:TYR:HE2	14	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1353)	1:8:A:GLY:H	1:37:A:CYS:H	10	0.11
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	4	0.11
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	4	0.11
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	4	0.11
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	8	0.11
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	8	0.11
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	8	0.11
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	20	0.11
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	20	0.11
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	20	0.11
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	7	0.11
(1,982)	1:42:A:GLN:H	1:42:A:GLN:HE22	1	0.11
(1,982)	1:42:A:GLN:H	1:42:A:GLN:HE22	2	0.11
(1,982)	1:42:A:GLN:H	1:42:A:GLN:HE22	14	0.11
(1,868)	1:102:A:ILE:HG21	1:121:A:ASP:H	11	0.11
(1,868)	1:102:A:ILE:HG22	1:121:A:ASP:H	11	0.11
(1,868)	1:102:A:ILE:HG23	1:121:A:ASP:H	11	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	2	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	2	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	2	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	7	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	7	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	7	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	9	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	9	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	9	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	10	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	10	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	10	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	11	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	11	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	11	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	12	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	12	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	12	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	15	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	15	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	15	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	17	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	17	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	17	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG11	18	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG12	18	0.11
(1,843)	1:100:A:VAL:H	1:100:A:VAL:HG13	18	0.11
(1,839)	1:91:A:THR:HG21	1:92:A:LEU:H	3	0.11
(1,839)	1:91:A:THR:HG22	1:92:A:LEU:H	3	0.11
(1,839)	1:91:A:THR:HG23	1:92:A:LEU:H	3	0.11
(1,839)	1:91:A:THR:HG21	1:92:A:LEU:H	6	0.11
(1,839)	1:91:A:THR:HG22	1:92:A:LEU:H	6	0.11
(1,839)	1:91:A:THR:HG23	1:92:A:LEU:H	6	0.11
(1,722)	1:95:A:ALA:HA	1:96:A:LEU:H	7	0.11
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG2	4	0.11
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG3	4	0.11
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG2	8	0.11
(1,594)	1:53:A:GLN:H	1:53:A:GLN:HG3	8	0.11
(1,580)	1:31:A:VAL:HB	1:34:A:SER:H	5	0.11
(1,566)	1:93:A:ARG:HD3	1:94:A:GLU:H	3	0.11
(1,566)	1:93:A:ARG:HD3	1:94:A:GLU:H	6	0.11
(1,566)	1:93:A:ARG:HD3	1:94:A:GLU:H	10	0.11
(1,496)	1:44:A:GLN:HA	1:44:A:GLN:HE22	10	0.11
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG21	8	0.11
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG22	8	0.11
(1,462)	1:96:A:LEU:HA	1:100:A:VAL:HG23	8	0.11
(1,455)	1:40:A:VAL:HG21	1:44:A:GLN:HB3	13	0.11
(1,455)	1:40:A:VAL:HG22	1:44:A:GLN:HB3	13	0.11
(1,455)	1:40:A:VAL:HG23	1:44:A:GLN:HB3	13	0.11
(1,415)	1:97:A:PRO:HB2	1:100:A:VAL:HG21	5	0.11
(1,415)	1:97:A:PRO:HB2	1:100:A:VAL:HG22	5	0.11
(1,415)	1:97:A:PRO:HB2	1:100:A:VAL:HG23	5	0.11
(1,409)	1:69:A:LYS:HG2	1:69:A:LYS:HD2	17	0.11
(1,409)	1:69:A:LYS:HG2	1:69:A:LYS:HD3	17	0.11
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD21	3	0.11
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD22	3	0.11
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD23	3	0.11
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD21	4	0.11
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD22	4	0.11
(1,396)	1:92:A:LEU:HA	1:92:A:LEU:HD23	4	0.11
(1,302)	1:40:A:VAL:HA	1:44:A:GLN:HE21	9	0.11
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD11	14	0.11
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD12	14	0.11
(1,88)	1:93:A:ARG:H	1:102:A:ILE:HD13	14	0.11
(1,69)	1:19:A:TYR:HB2	1:22:A:GLY:H	20	0.11
(2,5)	1:79:A:GLN:H	1:56:A:GLY:O	6	0.1
(1,2149)	1:93:A:ARG:HB2	1:94:A:GLU:H	1	0.1

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2149)	1:93:A:ARG:HB3	1:94:A:GLU:H	1	0.1
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD11	20	0.1
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD12	20	0.1
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD13	20	0.1
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD21	20	0.1
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD22	20	0.1
(1,2130)	1:86:A:GLN:H	1:87:A:LEU:HD23	20	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	3	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	3	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	3	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	3	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	3	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	3	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	9	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	9	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	9	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	9	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	9	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	9	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG21	13	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG22	13	0.1
(1,2109)	1:79:A:GLN:HE21	1:102:A:ILE:HG23	13	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG21	13	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG22	13	0.1
(1,2109)	1:79:A:GLN:HE22	1:102:A:ILE:HG23	13	0.1
(1,1877)	1:13:A:GLN:HE21	1:14:A:ASP:HB2	17	0.1
(1,1877)	1:13:A:GLN:HE21	1:14:A:ASP:HB3	17	0.1
(1,1877)	1:13:A:GLN:HE22	1:14:A:ASP:HB2	17	0.1
(1,1877)	1:13:A:GLN:HE22	1:14:A:ASP:HB3	17	0.1
(1,1858)	1:10:A:THR:H	1:14:A:ASP:HB2	9	0.1
(1,1858)	1:10:A:THR:H	1:14:A:ASP:HB3	9	0.1
(1,1762)	1:29:A:ILE:HA	1:57:A:VAL:HB	7	0.1
(1,1641)	1:41:A:GLU:HA	1:41:A:GLU:HG2	4	0.1
(1,1611)	1:13:A:GLN:HA	1:13:A:GLN:HG2	2	0.1
(1,1611)	1:13:A:GLN:HA	1:13:A:GLN:HG3	2	0.1
(1,1556)	1:79:A:GLN:H	1:79:A:GLN:HB3	5	0.1
(1,1502)	1:127:A:ASN:H	1:127:A:ASN:HD22	5	0.1
(1,1476)	1:29:A:ILE:HG21	1:56:A:GLY:HA2	7	0.1
(1,1476)	1:29:A:ILE:HG22	1:56:A:GLY:HA2	7	0.1
(1,1476)	1:29:A:ILE:HG23	1:56:A:GLY:HA2	7	0.1
(1,1353)	1:8:A:GLY:H	1:37:A:CYS:H	14	0.1
(1,1298)	1:75:A:LEU:H	1:75:A:LEU:HG	4	0.1

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	10	0.1
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	10	0.1
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	10	0.1
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	11	0.1
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	11	0.1
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	11	0.1
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	14	0.1
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	14	0.1
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	14	0.1
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	15	0.1
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	15	0.1
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	15	0.1
(1,1213)	1:29:A:ILE:HG21	1:30:A:PHE:H	18	0.1
(1,1213)	1:29:A:ILE:HG22	1:30:A:PHE:H	18	0.1
(1,1213)	1:29:A:ILE:HG23	1:30:A:PHE:H	18	0.1
(1,1208)	1:44:A:GLN:H	1:44:A:GLN:HB2	20	0.1
(1,1006)	1:60:A:ASN:H	1:61:A:HIS:H	19	0.1
(1,956)	1:12:A:GLY:HA3	1:13:A:GLN:HE21	4	0.1
(1,956)	1:12:A:GLY:HA3	1:13:A:GLN:HE21	10	0.1
(1,868)	1:102:A:ILE:HG21	1:121:A:ASP:H	15	0.1
(1,868)	1:102:A:ILE:HG22	1:121:A:ASP:H	15	0.1
(1,868)	1:102:A:ILE:HG23	1:121:A:ASP:H	15	0.1
(1,839)	1:91:A:THR:HG21	1:92:A:LEU:H	2	0.1
(1,839)	1:91:A:THR:HG22	1:92:A:LEU:H	2	0.1
(1,839)	1:91:A:THR:HG23	1:92:A:LEU:H	2	0.1
(1,839)	1:91:A:THR:HG21	1:92:A:LEU:H	12	0.1
(1,839)	1:91:A:THR:HG22	1:92:A:LEU:H	12	0.1
(1,839)	1:91:A:THR:HG23	1:92:A:LEU:H	12	0.1
(1,496)	1:44:A:GLN:HA	1:44:A:GLN:HE22	14	0.1
(1,376)	1:28:A:LEU:HD11	1:54:A:TYR:HB3	11	0.1
(1,376)	1:28:A:LEU:HD12	1:54:A:TYR:HB3	11	0.1
(1,376)	1:28:A:LEU:HD13	1:54:A:TYR:HB3	11	0.1
(1,376)	1:28:A:LEU:HD11	1:54:A:TYR:HB3	13	0.1
(1,376)	1:28:A:LEU:HD12	1:54:A:TYR:HB3	13	0.1
(1,376)	1:28:A:LEU:HD13	1:54:A:TYR:HB3	13	0.1
(1,367)	1:93:A:ARG:HB3	1:93:A:ARG:HD2	7	0.1
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD11	14	0.1
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD12	14	0.1
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD13	14	0.1
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD21	14	0.1
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD22	14	0.1
(1,359)	1:44:A:GLN:HB2	1:73:A:LEU:HD23	14	0.1

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,333)	1:127:A:ASN:H	1:127:A:ASN:HD22	5	0.1
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG21	8	0.1
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG22	8	0.1
(1,297)	1:30:A:PHE:HE1	1:66:A:VAL:HG23	8	0.1
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG21	8	0.1
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG22	8	0.1
(1,297)	1:30:A:PHE:HE2	1:66:A:VAL:HG23	8	0.1
(1,69)	1:19:A:TYR:HB2	1:22:A:GLY:H	9	0.1
(1,69)	1:19:A:TYR:HB2	1:22:A:GLY:H	16	0.1

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

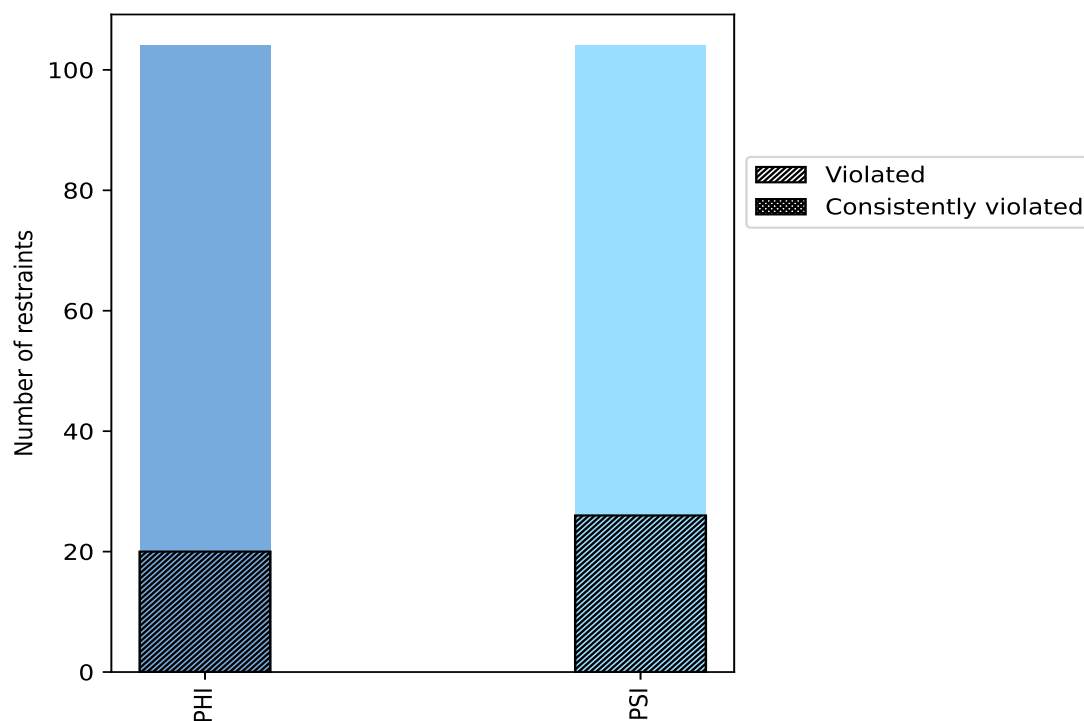
### 10.1 Summary of dihedral-angle violations [i](#)

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	% <sup>1</sup>	Violated <sup>3</sup>			Consistently Violated <sup>4</sup>		
			Count	% <sup>2</sup>	% <sup>1</sup>	Count	% <sup>2</sup>	% <sup>1</sup>
PHI	104	50.0	20	19.2	9.6	0	0.0	0.0
PSI	104	50.0	26	25.0	12.5	0	0.0	0.0
Total	208	100.0	46	22.1	22.1	0	0.0	0.0

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

#### 10.1.1 Bar chart : Distribution of dihedral-angles and violations [i](#)



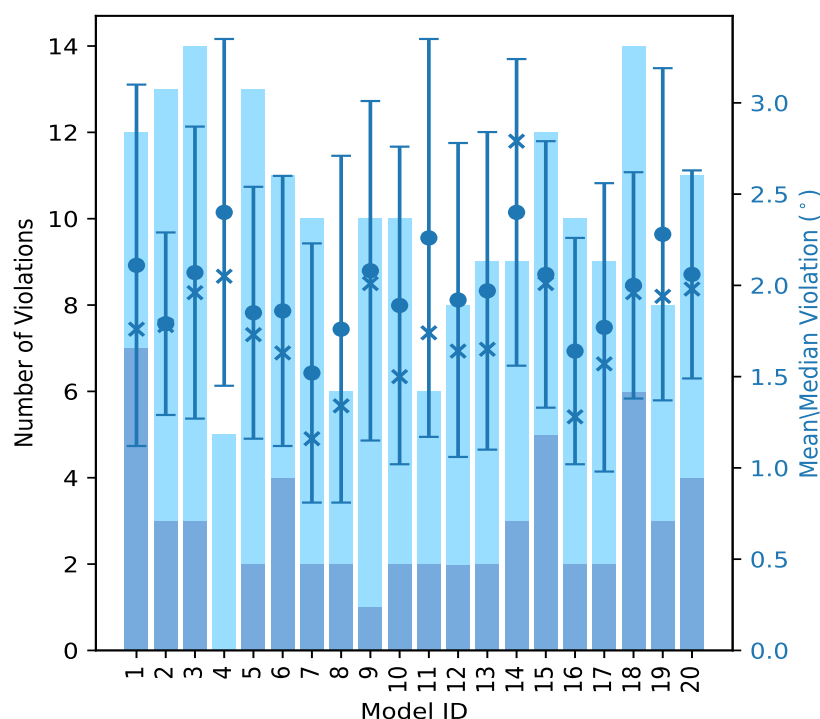
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

## 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	7	5	12	2.11	4.4	0.99	1.76
2	3	10	13	1.79	2.87	0.5	1.78
3	3	11	14	2.07	4.16	0.8	1.96
4	0	5	5	2.4	4.24	0.95	2.05
5	2	11	13	1.85	3.54	0.69	1.73
6	4	7	11	1.86	3.42	0.74	1.63
7	2	8	10	1.52	3.15	0.71	1.16
8	2	4	6	1.76	3.8	0.95	1.34
9	1	9	10	2.08	4.44	0.93	2.01
10	2	8	10	1.89	4.03	0.87	1.5
11	2	4	6	2.26	4.26	1.09	1.74
12	2	6	8	1.92	3.71	0.86	1.64
13	2	7	9	1.97	3.95	0.87	1.65
14	3	6	9	2.4	3.82	0.84	2.79
15	5	7	12	2.06	3.33	0.73	2.01
16	2	8	10	1.64	2.56	0.62	1.28
17	2	7	9	1.77	3.95	0.79	1.57
18	6	8	14	2.0	3.21	0.62	1.96
19	3	5	8	2.28	3.8	0.91	1.94
20	4	7	11	2.06	3.06	0.57	1.98

### 10.2.1 Bar graph : Dihedral violation statistics for each model [i](#)



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

### 10.3 Dihedral-angle violation statistics for the ensemble [i](#)

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

Number of violated restraints			Fraction of the ensemble	
PHI	PSI	Total	Count <sup>1</sup>	%
7	6	13	1	5.0
2	2	4	2	10.0
6	6	12	3	15.0
3	2	5	4	20.0
1	0	1	5	25.0
0	2	2	6	30.0
0	1	1	7	35.0
0	3	3	8	40.0
0	0	0	9	45.0
0	0	0	10	50.0
1	0	1	11	55.0

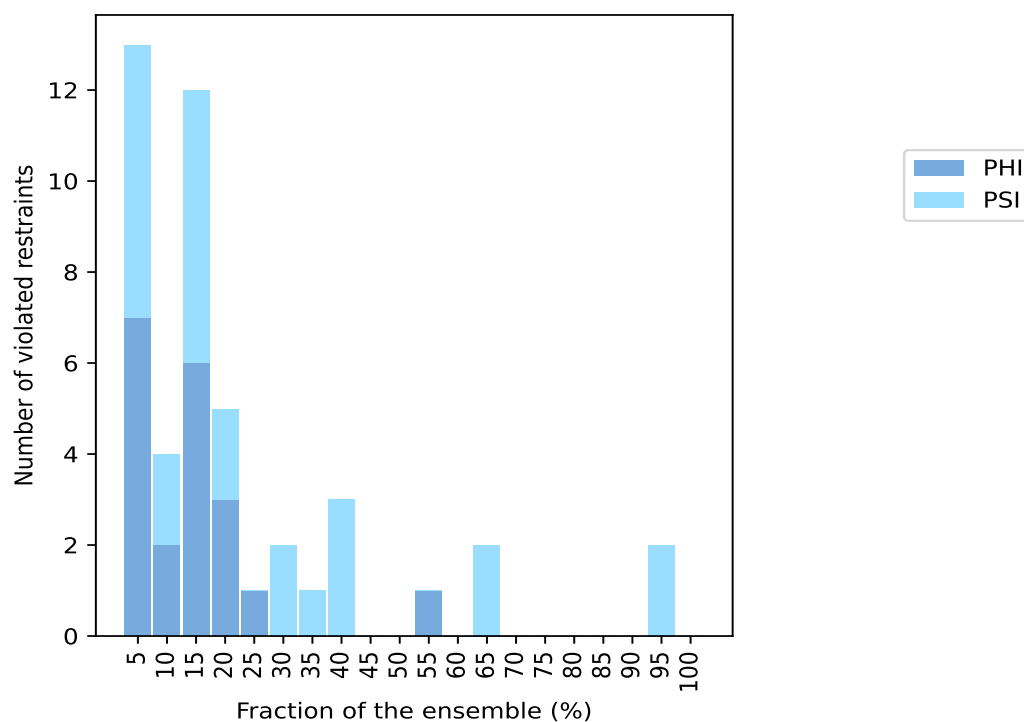
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Number of violated restraints			Fraction of the ensemble	
PHI	PSI	Total	Count <sup>1</sup>	%
0	0	0	12	60.0
0	2	2	13	65.0
0	0	0	14	70.0
0	0	0	15	75.0
0	0	0	16	80.0
0	0	0	17	85.0
0	0	0	18	90.0
0	2	2	19	95.0
0	0	0	20	100.0

<sup>1</sup> Number of models with violations

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

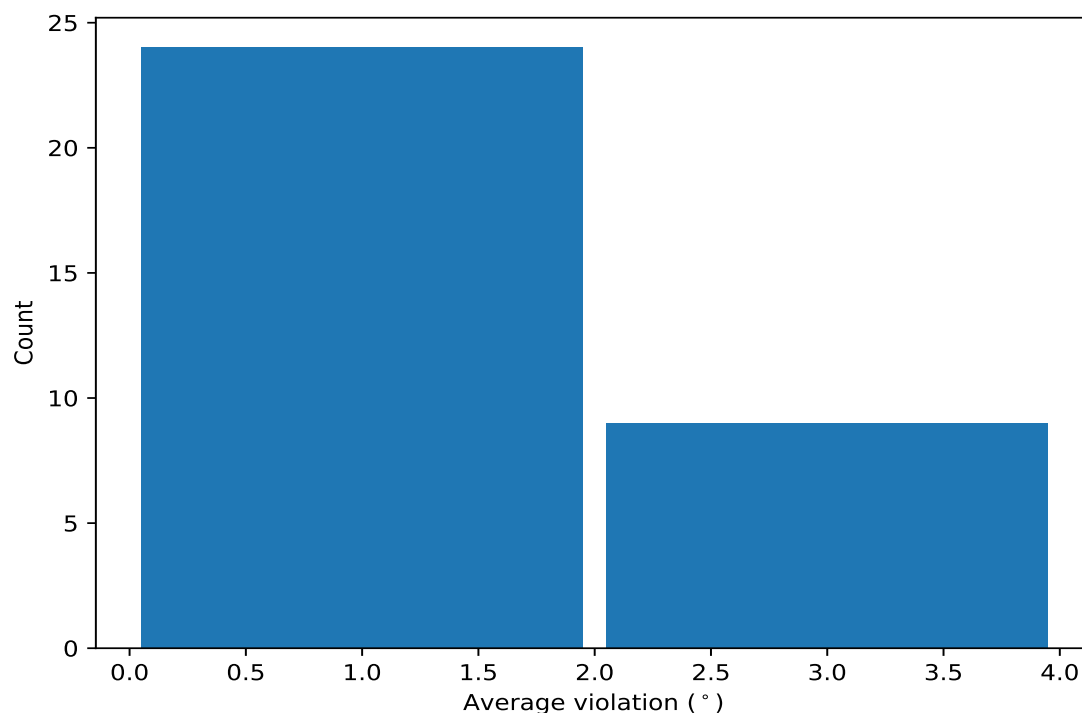


## 10.4 Most violated dihedral-angle restraints in the ensemble ⓘ

### 10.4.1 Histogram : Distribution of mean dihedral-angle violations ⓘ

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

in the ensemble



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

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

Key	Atom-1	Atom-2	Atom-3	Atom-4	Models <sup>1</sup>	Mean	SD <sup>2</sup>	Median
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	19	2.96	0.81	3.1
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	19	2.33	0.76	2.06
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	13	2.16	0.62	2.13
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	13	1.76	0.42	1.68
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	11	2.24	0.63	2.04
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	8	2.12	1.04	1.76
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	8	2.06	0.78	2.12
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	8	1.58	0.36	1.61
(1,34)	1:24:A:ILE:N	1:24:A:ILE:CA	1:24:A:ILE:C	1:25:A:TYR:N	7	2.11	0.6	2.15
(1,132)	1:82:A:GLY:N	1:82:A:GLY:CA	1:82:A:GLY:C	1:83:A:ASN:N	6	2.06	1.09	1.54
(1,94)	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	1:60:A:ASN:N	6	1.73	0.64	1.7
(1,5)	1:6:A:VAL:C	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	5	1.87	1.28	1.28
(1,129)	1:79:A:GLN:C	1:80:A:LEU:N	1:80:A:LEU:CA	1:80:A:LEU:C	4	2.68	0.61	2.86
(1,133)	1:82:A:GLY:C	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	4	1.7	0.48	1.62
(1,140)	1:87:A:LEU:N	1:87:A:LEU:CA	1:87:A:LEU:C	1:88:A:TYR:N	4	1.47	0.24	1.44
(1,160)	1:97:A:PRO:N	1:97:A:PRO:CA	1:97:A:PRO:C	1:98:A:ALA:N	4	1.18	0.07	1.16
(1,43)	1:28:A:LEU:C	1:29:A:ILE:N	1:29:A:ILE:CA	1:29:A:ILE:C	4	1.16	0.19	1.07
(1,178)	1:106:A:LEU:N	1:106:A:LEU:CA	1:106:A:LEU:C	1:107:A:SER:N	3	1.77	0.1	1.78
(1,135)	1:83:A:ASN:C	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	3	1.77	0.35	1.72
(1,95)	1:59:A:ARG:C	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	3	1.69	0.63	1.36

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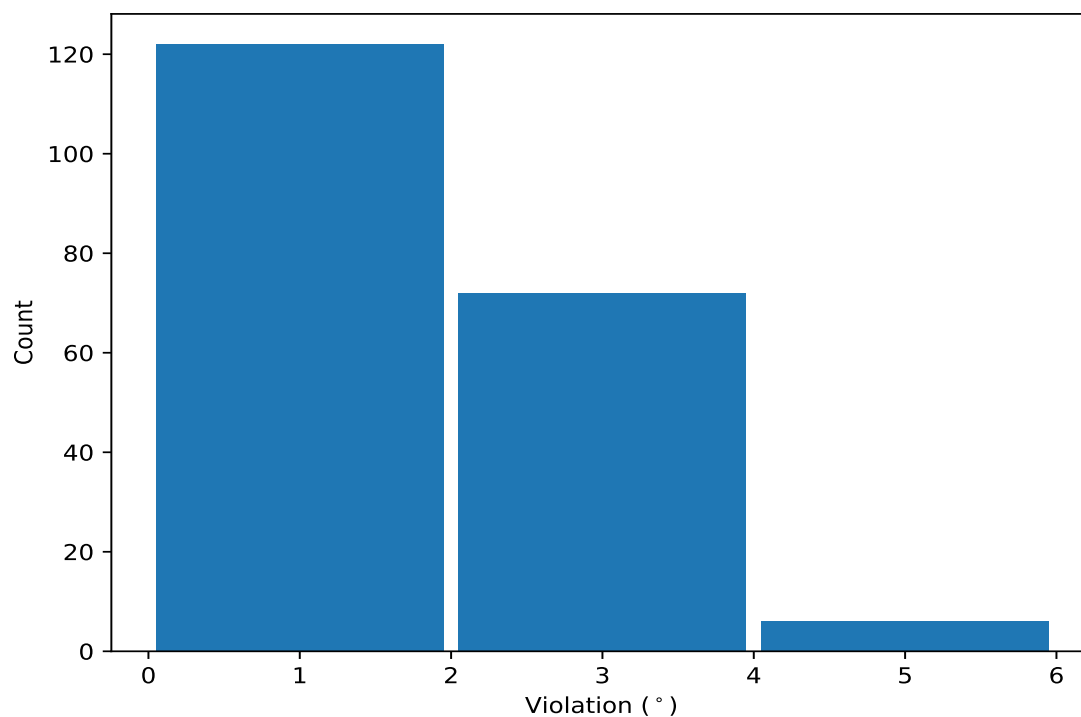
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models <sup>1</sup>	Mean	SD <sup>2</sup>	Median
(1,182)	1:108:A:VAL:N	1:108:A:VAL:CA	1:108:A:VAL:C	1:109:A:GLY:N	3	1.6	0.2	1.71
(1,32)	1:21:A:ALA:N	1:21:A:ALA:CA	1:21:A:ALA:C	1:22:A:GLY:N	3	1.58	0.48	1.29
(1,137)	1:85:A:GLU:C	1:86:A:GLN:N	1:86:A:GLN:CA	1:86:A:GLN:C	3	1.44	0.44	1.22
(1,186)	1:113:A:PRO:N	1:113:A:PRO:CA	1:113:A:PRO:C	1:114:A:ALA:N	3	1.43	0.27	1.48
(1,180)	1:107:A:SER:N	1:107:A:SER:CA	1:107:A:SER:C	1:108:A:VAL:N	3	1.4	0.23	1.26
(1,147)	1:90:A:ASP:C	1:91:A:THR:N	1:91:A:THR:CA	1:91:A:THR:C	3	1.37	0.19	1.32
(1,173)	1:103:A:TRP:C	1:104:A:LYS:N	1:104:A:LYS:CA	1:104:A:LYS:C	3	1.37	0.17	1.4
(1,164)	1:99:A:HIS:N	1:99:A:HIS:CA	1:99:A:HIS:C	1:100:A:VAL:N	3	1.25	0.22	1.17
(1,121)	1:75:A:LEU:C	1:76:A:ALA:N	1:76:A:ALA:CA	1:76:A:ALA:C	3	1.07	0.02	1.06
(1,8)	1:8:A:GLY:N	1:8:A:GLY:CA	1:8:A:GLY:C	1:9:A:LEU:N	2	1.62	0.58	1.62
(1,16)	1:13:A:GLN:N	1:13:A:GLN:CA	1:13:A:GLN:C	1:14:A:ASP:N	2	1.57	0.5	1.57
(1,55)	1:37:A:CYS:C	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	2	1.52	0.18	1.52
(1,93)	1:58:A:PHE:C	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	2	1.23	0.09	1.23

<sup>1</sup> Number of violated models, <sup>2</sup>Standard 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 (°)
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	9	4.44
(1,5)	1:6:A:VAL:C	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1	4.4
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	11	4.26
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	4	4.24
(1,132)	1:82:A:GLY:N	1:82:A:GLY:CA	1:82:A:GLY:C	1:83:A:ASN:N	3	4.16
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	10	4.03
(1,7)	1:7:A:CYS:C	1:8:A:GLY:N	1:8:A:GLY:CA	1:8:A:GLY:C	13	3.95
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	17	3.95
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	14	3.82
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	8	3.8
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	19	3.8
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	12	3.71
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	5	3.54
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	1	3.43
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	6	3.42
(1,129)	1:79:A:GLN:C	1:80:A:LEU:N	1:80:A:LEU:CA	1:80:A:LEU:C	15	3.33
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	19	3.3
(1,4)	1:6:A:VAL:N	1:6:A:VAL:CA	1:6:A:VAL:C	1:7:A:CYS:N	1	3.24
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	18	3.21
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	11	3.15
(1,90)	1:57:A:VAL:N	1:57:A:VAL:CA	1:57:A:VAL:C	1:58:A:PHE:N	7	3.15
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	15	3.1
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	20	3.06
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	14	2.98
(1,34)	1:24:A:ILE:N	1:24:A:ILE:CA	1:24:A:ILE:C	1:25:A:TYR:N	19	2.98
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	14	2.92
(1,129)	1:79:A:GLN:C	1:80:A:LEU:N	1:80:A:LEU:CA	1:80:A:LEU:C	14	2.92
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	13	2.91
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	3	2.91
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	2	2.87
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	12	2.84
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	20	2.8
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	15	2.8
(1,132)	1:82:A:GLY:N	1:82:A:GLY:CA	1:82:A:GLY:C	1:83:A:ASN:N	14	2.79
(1,129)	1:79:A:GLN:C	1:80:A:LEU:N	1:80:A:LEU:CA	1:80:A:LEU:C	18	2.79
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	9	2.68
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	10	2.63
(1,48)	1:32:A:ALA:N	1:32:A:ALA:CA	1:32:A:ALA:C	1:33:A:THR:N	5	2.62
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	18	2.61
(1,95)	1:59:A:ARG:C	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	3	2.57
(1,34)	1:24:A:ILE:N	1:24:A:ILE:CA	1:24:A:ILE:C	1:25:A:TYR:N	16	2.56
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	16	2.56
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	2	2.54
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	10	2.53
(1,94)	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	1:60:A:ASN:N	20	2.52
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	6	2.5
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	7	2.5

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,131)	1:81:A:HIS:C	1:82:A:GLY:N	1:82:A:GLY:CA	1:82:A:GLY:C	18	2.48
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	16	2.48
(1,34)	1:24:A:ILE:N	1:24:A:ILE:CA	1:24:A:ILE:C	1:25:A:TYR:N	20	2.46
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	3	2.46
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	15	2.4
(1,133)	1:82:A:GLY:C	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	3	2.4
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	6	2.39
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	15	2.36
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	9	2.35
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	6	2.35
(1,94)	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	1:60:A:ASN:N	3	2.34
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	18	2.32
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	5	2.3
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	6	2.29
(1,181)	1:107:A:SER:C	1:108:A:VAL:N	1:108:A:VAL:CA	1:108:A:VAL:C	5	2.27
(1,32)	1:21:A:ALA:N	1:21:A:ALA:CA	1:21:A:ALA:C	1:22:A:GLY:N	4	2.25
(1,135)	1:83:A:ASN:C	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	20	2.22
(1,94)	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	1:60:A:ASN:N	18	2.2
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	2	2.2
(1,8)	1:8:A:GLY:N	1:8:A:GLY:CA	1:8:A:GLY:C	1:9:A:LEU:N	13	2.2
(1,34)	1:24:A:ILE:N	1:24:A:ILE:CA	1:24:A:ILE:C	1:25:A:TYR:N	9	2.15
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	15	2.13
(1,16)	1:13:A:GLN:N	1:13:A:GLN:CA	1:13:A:GLN:C	1:14:A:ASP:N	9	2.07
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	18	2.06
(1,137)	1:85:A:GLU:C	1:86:A:GLN:N	1:86:A:GLN:CA	1:86:A:GLN:C	2	2.05
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	4	2.05
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	5	2.04
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	12	2.04
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	1	2.04
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	3	2.01
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	14	2.01
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	19	1.98
(1,34)	1:24:A:ILE:N	1:24:A:ILE:CA	1:24:A:ILE:C	1:25:A:TYR:N	2	1.98
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	20	1.98
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	9	1.96
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	3	1.92
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	19	1.91
(1,178)	1:106:A:LEU:N	1:106:A:LEU:CA	1:106:A:LEU:C	1:107:A:SER:N	4	1.89
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	7	1.89
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	15	1.89
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	19	1.87
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	3	1.86
(1,133)	1:82:A:GLY:C	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	18	1.85
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	1	1.84
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	8	1.83
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	11	1.82
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	17	1.82
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	16	1.82
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	5	1.81
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	2	1.8
(1,140)	1:87:A:LEU:N	1:87:A:LEU:CA	1:87:A:LEU:C	1:88:A:TYR:N	18	1.8

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	20	1.79
(1,178)	1:106:A:LEU:N	1:106:A:LEU:CA	1:106:A:LEU:C	1:107:A:SER:N	18	1.78
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	2	1.78
(1,182)	1:108:A:VAL:N	1:108:A:VAL:CA	1:108:A:VAL:C	1:109:A:GLY:N	1	1.77
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	1	1.75
(1,186)	1:113:A:PRO:N	1:113:A:PRO:CA	1:113:A:PRO:C	1:114:A:ALA:N	5	1.73
(1,180)	1:107:A:SER:N	1:107:A:SER:CA	1:107:A:SER:C	1:108:A:VAL:N	5	1.72
(1,135)	1:83:A:ASN:C	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	14	1.72
(1,182)	1:108:A:VAL:N	1:108:A:VAL:CA	1:108:A:VAL:C	1:109:A:GLY:N	13	1.71
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	17	1.71
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	12	1.71
(1,55)	1:37:A:CYS:C	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1	1.7
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	15	1.68
(1,129)	1:79:A:GLN:C	1:80:A:LEU:N	1:80:A:LEU:CA	1:80:A:LEU:C	20	1.68
(1,178)	1:106:A:LEU:N	1:106:A:LEU:CA	1:106:A:LEU:C	1:107:A:SER:N	11	1.65
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	17	1.65
(1,6)	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	1:8:A:GLY:N	13	1.65
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	6	1.63
(1,147)	1:90:A:ASP:C	1:91:A:THR:N	1:91:A:THR:CA	1:91:A:THR:C	15	1.63
(1,140)	1:87:A:LEU:N	1:87:A:LEU:CA	1:87:A:LEU:C	1:88:A:TYR:N	10	1.6
(1,34)	1:24:A:ILE:N	1:24:A:ILE:CA	1:24:A:ILE:C	1:25:A:TYR:N	12	1.58
(1,132)	1:82:A:GLY:N	1:82:A:GLY:CA	1:82:A:GLY:C	1:83:A:ASN:N	17	1.57
(1,173)	1:103:A:TRP:C	1:104:A:LYS:N	1:104:A:LYS:CA	1:104:A:LYS:C	11	1.56
(1,164)	1:99:A:HIS:N	1:99:A:HIS:CA	1:99:A:HIS:C	1:100:A:VAL:N	4	1.55
(1,5)	1:6:A:VAL:C	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	10	1.55
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	13	1.53
(1,46)	1:31:A:VAL:N	1:31:A:VAL:CA	1:31:A:VAL:C	1:32:A:ALA:N	5	1.53
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	3	1.52
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	20	1.52
(1,132)	1:82:A:GLY:N	1:82:A:GLY:CA	1:82:A:GLY:C	1:83:A:ASN:N	2	1.52
(1,165)	1:99:A:HIS:C	1:100:A:VAL:N	1:100:A:VAL:CA	1:100:A:VAL:C	1	1.51
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	9	1.51
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	20	1.5
(1,186)	1:113:A:PRO:N	1:113:A:PRO:CA	1:113:A:PRO:C	1:114:A:ALA:N	2	1.48
(1,57)	1:38:A:VAL:C	1:39:A:ASN:N	1:39:A:ASN:CA	1:39:A:ASN:C	8	1.48
(1,43)	1:28:A:LEU:C	1:29:A:ILE:N	1:29:A:ILE:CA	1:29:A:ILE:C	6	1.48
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	10	1.44
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	13	1.44
(1,35)	1:24:A:ILE:C	1:25:A:TYR:N	1:25:A:TYR:CA	1:25:A:TYR:C	3	1.41
(1,173)	1:103:A:TRP:C	1:104:A:LYS:N	1:104:A:LYS:CA	1:104:A:LYS:C	19	1.4
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	2	1.4
(1,133)	1:82:A:GLY:C	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	15	1.39
(1,135)	1:83:A:ASN:C	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	18	1.37
(1,192)	1:117:A:PHE:N	1:117:A:PHE:CA	1:117:A:PHE:C	1:118:A:GLN:N	14	1.36
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	10	1.36
(1,95)	1:59:A:ARG:C	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1	1.36
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	10	1.34
(1,55)	1:37:A:CYS:C	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	13	1.33
(1,147)	1:90:A:ASP:C	1:91:A:THR:N	1:91:A:THR:CA	1:91:A:THR:C	17	1.32
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	17	1.31
(1,182)	1:108:A:VAL:N	1:108:A:VAL:CA	1:108:A:VAL:C	1:109:A:GLY:N	2	1.31

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,93)	1:58:A:PHE:C	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	5	1.31
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	17	1.3
(1,160)	1:97:A:PRO:N	1:97:A:PRO:CA	1:97:A:PRO:C	1:98:A:ALA:N	17	1.29
(1,32)	1:21:A:ALA:N	1:21:A:ALA:CA	1:21:A:ALA:C	1:22:A:GLY:N	16	1.29
(1,5)	1:6:A:VAL:C	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	12	1.28
(1,140)	1:87:A:LEU:N	1:87:A:LEU:CA	1:87:A:LEU:C	1:88:A:TYR:N	16	1.27
(1,180)	1:107:A:SER:N	1:107:A:SER:CA	1:107:A:SER:C	1:108:A:VAL:N	2	1.26
(1,132)	1:82:A:GLY:N	1:82:A:GLY:CA	1:82:A:GLY:C	1:83:A:ASN:N	18	1.26
(1,56)	1:38:A:VAL:N	1:38:A:VAL:CA	1:38:A:VAL:C	1:39:A:ASN:N	9	1.26
(1,140)	1:87:A:LEU:N	1:87:A:LEU:CA	1:87:A:LEU:C	1:88:A:TYR:N	7	1.22
(1,137)	1:85:A:GLU:C	1:86:A:GLN:N	1:86:A:GLN:CA	1:86:A:GLN:C	7	1.22
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	10	1.22
(1,180)	1:107:A:SER:N	1:107:A:SER:CA	1:107:A:SER:C	1:108:A:VAL:N	8	1.21
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	9	1.21
(1,94)	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	1:60:A:ASN:N	16	1.2
(1,32)	1:21:A:ALA:N	1:21:A:ALA:CA	1:21:A:ALA:C	1:22:A:GLY:N	3	1.2
(1,164)	1:99:A:HIS:N	1:99:A:HIS:CA	1:99:A:HIS:C	1:100:A:VAL:N	3	1.17
(1,160)	1:97:A:PRO:N	1:97:A:PRO:CA	1:97:A:PRO:C	1:98:A:ALA:N	5	1.17
(1,147)	1:90:A:ASP:C	1:91:A:THR:N	1:91:A:THR:CA	1:91:A:THR:C	1	1.17
(1,200)	1:122:A:LYS:N	1:122:A:LYS:CA	1:122:A:LYS:C	1:123:A:TYR:N	10	1.16
(1,160)	1:97:A:PRO:N	1:97:A:PRO:CA	1:97:A:PRO:C	1:98:A:ALA:N	16	1.15
(1,134)	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	1:84:A:GLU:N	8	1.15
(1,133)	1:82:A:GLY:C	1:83:A:ASN:N	1:83:A:ASN:CA	1:83:A:ASN:C	20	1.15
(1,173)	1:103:A:TRP:C	1:104:A:LYS:N	1:104:A:LYS:CA	1:104:A:LYS:C	6	1.14
(1,95)	1:59:A:ARG:C	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	18	1.14
(1,93)	1:58:A:PHE:C	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	11	1.14
(1,9)	1:9:A:LEU:C	1:10:A:THR:N	1:10:A:THR:CA	1:10:A:THR:C	12	1.14
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	6	1.12
(1,10)	1:10:A:THR:N	1:10:A:THR:CA	1:10:A:THR:C	1:11:A:ARG:N	9	1.12
(1,157)	1:95:A:ALA:C	1:96:A:LEU:N	1:96:A:LEU:CA	1:96:A:LEU:C	14	1.11
(1,160)	1:97:A:PRO:N	1:97:A:PRO:CA	1:97:A:PRO:C	1:98:A:ALA:N	6	1.1
(1,121)	1:75:A:LEU:C	1:76:A:ALA:N	1:76:A:ALA:CA	1:76:A:ALA:C	2	1.1
(1,5)	1:6:A:VAL:C	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	18	1.1
(1,43)	1:28:A:LEU:C	1:29:A:ILE:N	1:29:A:ILE:CA	1:29:A:ILE:C	7	1.09
(1,186)	1:113:A:PRO:N	1:113:A:PRO:CA	1:113:A:PRO:C	1:114:A:ALA:N	7	1.08
(1,94)	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	1:60:A:ASN:N	12	1.07
(1,16)	1:13:A:GLN:N	1:13:A:GLN:CA	1:13:A:GLN:C	1:14:A:ASP:N	3	1.07
(1,121)	1:75:A:LEU:C	1:76:A:ALA:N	1:76:A:ALA:CA	1:76:A:ALA:C	8	1.06
(1,132)	1:82:A:GLY:N	1:82:A:GLY:CA	1:82:A:GLY:C	1:83:A:ASN:N	7	1.05
(1,121)	1:75:A:LEU:C	1:76:A:ALA:N	1:76:A:ALA:CA	1:76:A:ALA:C	6	1.05
(1,43)	1:28:A:LEU:C	1:29:A:ILE:N	1:29:A:ILE:CA	1:29:A:ILE:C	1	1.05
(1,8)	1:8:A:GLY:N	1:8:A:GLY:CA	1:8:A:GLY:C	1:9:A:LEU:N	16	1.05
(1,137)	1:85:A:GLU:C	1:86:A:GLN:N	1:86:A:GLN:CA	1:86:A:GLN:C	15	1.04
(1,94)	1:59:A:ARG:N	1:59:A:ARG:CA	1:59:A:ARG:C	1:60:A:ASN:N	15	1.03
(1,34)	1:24:A:ILE:N	1:24:A:ILE:CA	1:24:A:ILE:C	1:25:A:TYR:N	5	1.03
(1,22)	1:16:A:LYS:N	1:16:A:LYS:CA	1:16:A:LYS:C	1:17:A:ALA:N	7	1.03
(1,164)	1:99:A:HIS:N	1:99:A:HIS:CA	1:99:A:HIS:C	1:100:A:VAL:N	13	1.02
(1,136)	1:84:A:GLU:N	1:84:A:GLU:CA	1:84:A:GLU:C	1:85:A:GLU:N	5	1.02
(1,43)	1:28:A:LEU:C	1:29:A:ILE:N	1:29:A:ILE:CA	1:29:A:ILE:C	16	1.02
(1,96)	1:60:A:ASN:N	1:60:A:ASN:CA	1:60:A:ASN:C	1:61:A:HIS:N	7	1.01
(1,5)	1:6:A:VAL:C	1:7:A:CYS:N	1:7:A:CYS:CA	1:7:A:CYS:C	19	1.01