



Full wwPDB NMR Structure Validation Report ⓘ

Mar 6, 2026 – 03:12 PM UTC

PDB ID : 2M9V / pdb_00002m9v
BMRB ID : 19311
Title : Structure of Saccharomyces cerevisiae Est3 protein
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Deposited on : 2013-06-21

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with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
BMRB Restraints Analysis : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

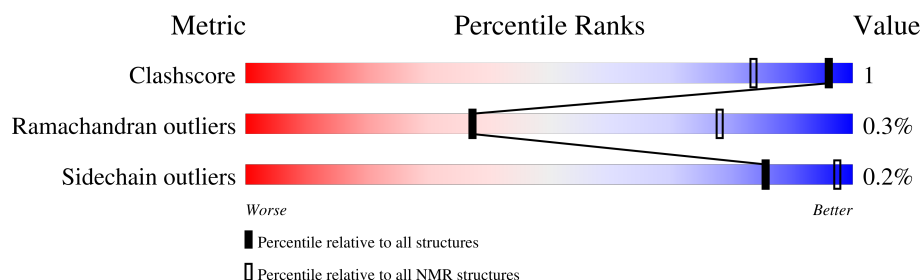
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 47%.

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



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

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

Mol	Chain	Length	Quality of chain
1	A	170	

2 Ensemble composition and analysis

This entry contains 20 models. Model 1 is the overall representative, medoid model (most similar to other models).

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

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:18-A:29, A:39-A:169, A:176-A:181 (149)	1.24	1

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

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

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

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2675 atoms, of which 1319 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Telomere replication protein EST3.

Mol	Chain	Residues	Atoms						Trace
1	A	170	Total	C	H	N	O	S	0
			2675	863	1319	224	261	8	

There are 2 discrepancies between the modelled and reference sequences:

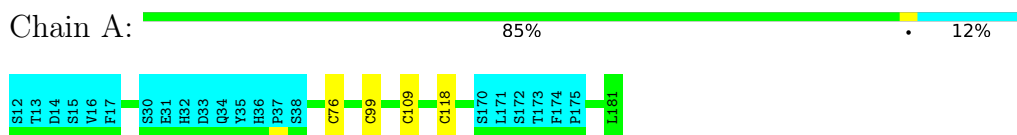
Chain	Residue	Modelled	Actual	Comment	Reference
A	12	SER	-	expression tag	UNP Q03096
A	142	SER	CYS	engineered mutation	UNP Q03096

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: Telomere replication protein EST3

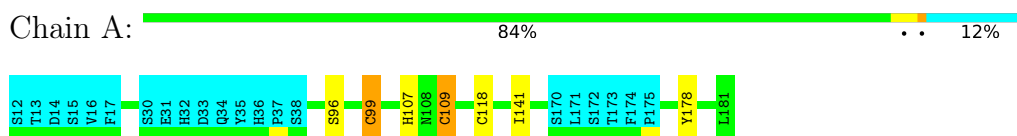


4.2 Scores per residue for each member of the ensemble

Colouring as in section [4.1](#) above.

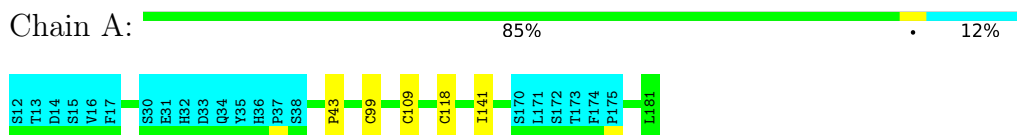
4.2.1 Score per residue for model 1 (medoid)

- Molecule 1: Telomere replication protein EST3



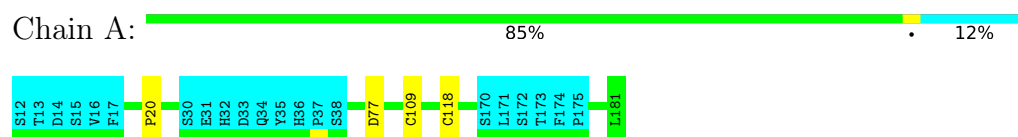
4.2.2 Score per residue for model 2

- Molecule 1: Telomere replication protein EST3



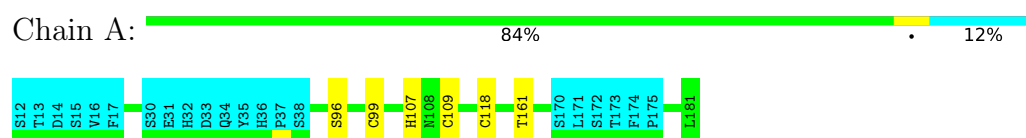
4.2.3 Score per residue for model 3

- Molecule 1: Telomere replication protein EST3



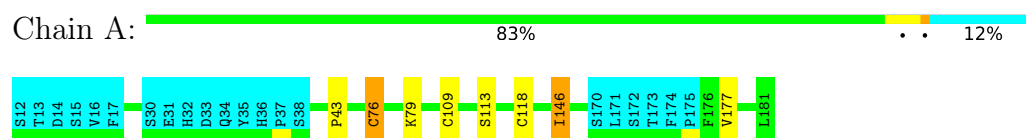
4.2.4 Score per residue for model 4

- Molecule 1: Telomere replication protein EST3



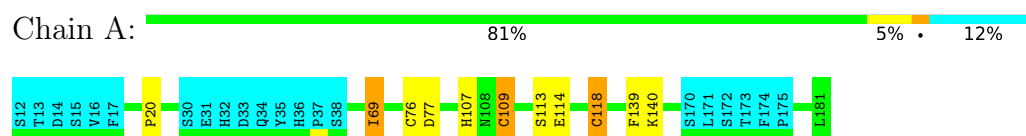
4.2.5 Score per residue for model 5

- Molecule 1: Telomere replication protein EST3



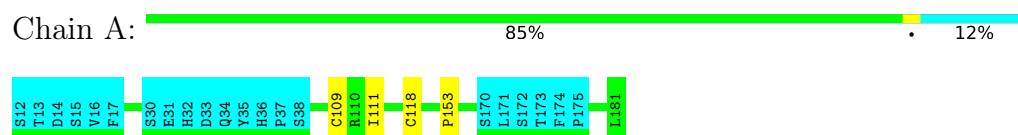
4.2.6 Score per residue for model 6

- Molecule 1: Telomere replication protein EST3



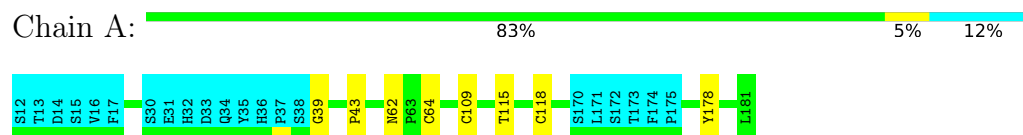
4.2.7 Score per residue for model 7

- Molecule 1: Telomere replication protein EST3



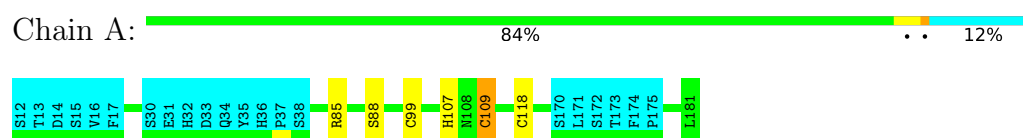
4.2.8 Score per residue for model 8

- Molecule 1: Telomere replication protein EST3



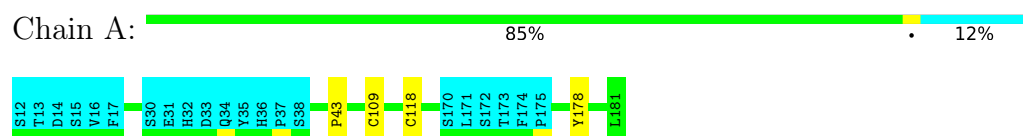
4.2.9 Score per residue for model 9

- Molecule 1: Telomere replication protein EST3



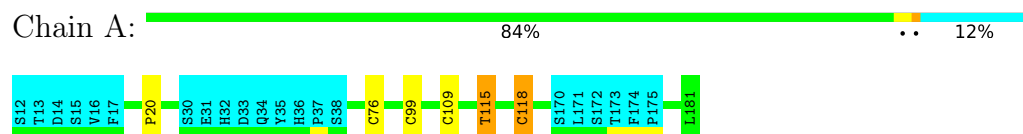
4.2.10 Score per residue for model 10

- Molecule 1: Telomere replication protein EST3



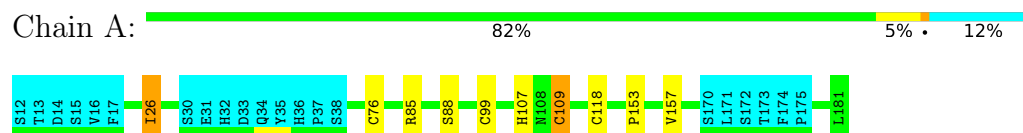
4.2.11 Score per residue for model 11

- Molecule 1: Telomere replication protein EST3



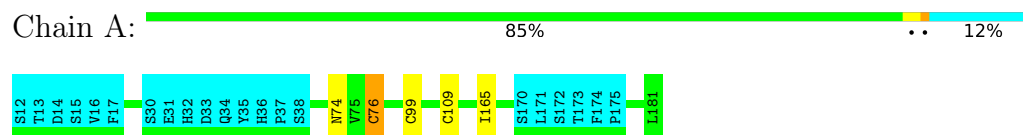
4.2.12 Score per residue for model 12

- Molecule 1: Telomere replication protein EST3



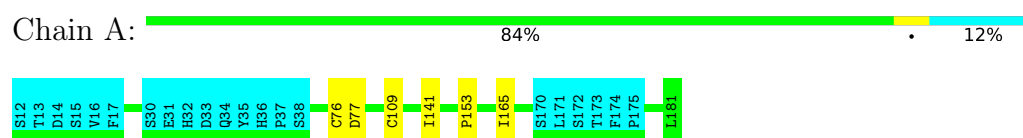
4.2.13 Score per residue for model 13

- Molecule 1: Telomere replication protein EST3



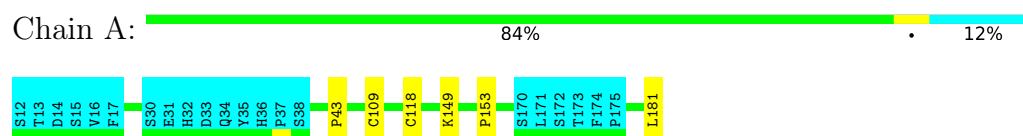
4.2.14 Score per residue for model 14

- Molecule 1: Telomere replication protein EST3



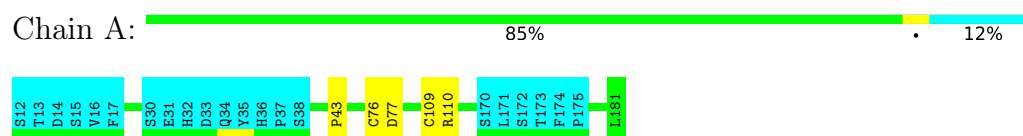
4.2.15 Score per residue for model 15

- Molecule 1: Telomere replication protein EST3



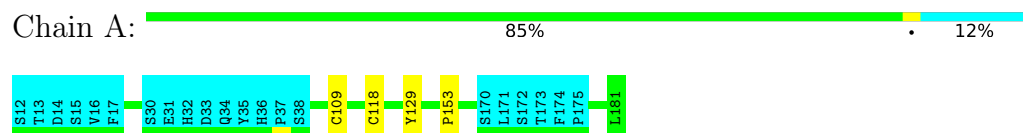
4.2.16 Score per residue for model 16

- Molecule 1: Telomere replication protein EST3



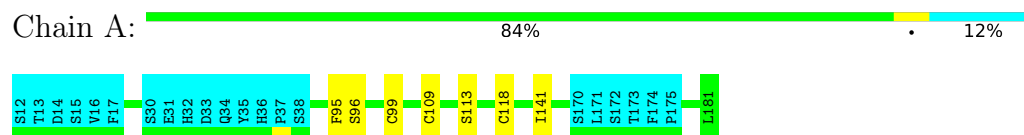
4.2.17 Score per residue for model 17

- Molecule 1: Telomere replication protein EST3



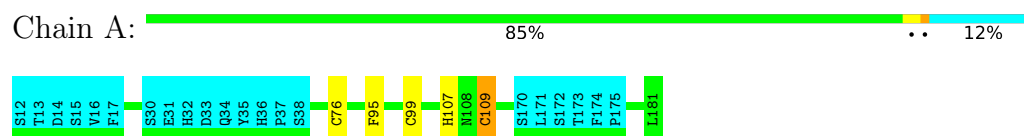
4.2.18 Score per residue for model 18

- Molecule 1: Telomere replication protein EST3



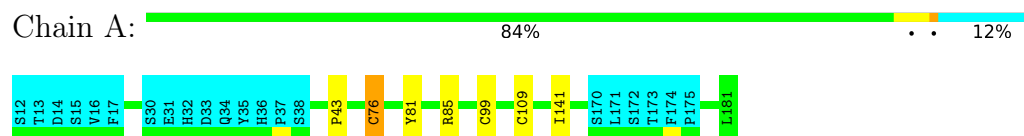
4.2.19 Score per residue for model 19

- Molecule 1: Telomere replication protein EST3



4.2.20 Score per residue for model 20

- Molecule 1: Telomere replication protein EST3



5 Refinement protocol and experimental data overview

The models were refined using the following method: *Molecular fragment replacement*.

Of the 500 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
Rosetta	structure solution	3.4
CS-Rosetta toolbox	structure solution	1.3
Rosetta	refinement	3.4

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	1071
Number of shifts mapped to atoms	1071
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	47%

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	0.56±0.01	1±1/1215 (0.1± 0.0%)	0.73±0.01	2±1/1651 (0.1± 0.0%)
All	All	0.56	15/24300 (0.1%)	0.73	49/33020 (0.1%)

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	43	PRO	N-CD	6.69	1.57	1.47	20	7
1	A	153	PRO	N-CD	5.41	1.55	1.47	17	5
1	A	20	PRO	N-CD	5.26	1.55	1.47	11	3

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	76	CYS	N-CA-CB	-5.68	102.75	111.56	11	9
1	A	109	CYS	N-CA-CB	-5.68	102.76	111.56	13	20
1	A	118	CYS	N-CA-CB	-5.57	102.70	111.05	17	14
1	A	99	CYS	N-CA-CB	-5.01	102.74	110.01	11	6

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	1189	1178	1184	2±2
All	All	23780	23560	23680	44

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:141:ILE:HG23	1:A:141:ILE:O	0.62	1.93	20	4
1:A:85:ARG:NH2	1:A:88:SER:O	0.61	2.33	12	1
1:A:69:ILE:HD12	1:A:69:ILE:N	0.60	2.12	6	1
1:A:26:ILE:C	1:A:26:ILE:HD12	0.56	2.26	12	1
1:A:111:ILE:O	1:A:111:ILE:HG22	0.54	2.03	7	1
1:A:107:HIS:HB3	1:A:109:CYS:SG	0.53	2.43	19	5
1:A:146:ILE:O	1:A:146:ILE:HG12	0.52	2.05	5	1
1:A:62:ASN:HB3	1:A:64:CYS:SG	0.50	2.47	8	1
1:A:76:CYS:HG	1:A:81:TYR:HE2	0.50	1.49	20	1
1:A:69:ILE:N	1:A:69:ILE:CD1	0.50	2.74	6	1
1:A:76:CYS:SG	1:A:81:TYR:CE2	0.47	3.08	20	1
1:A:95:PHE:HB3	1:A:99:CYS:SG	0.47	2.49	18	2
1:A:99:CYS:SG	1:A:157:VAL:O	0.46	2.74	12	1
1:A:96:SER:O	1:A:99:CYS:HB3	0.46	2.10	4	3
1:A:178:TYR:CD1	1:A:178:TYR:N	0.46	2.83	8	2
1:A:115:THR:HA	1:A:118:CYS:SG	0.46	2.50	11	2
1:A:76:CYS:SG	1:A:81:TYR:HE2	0.45	2.34	20	1
1:A:26:ILE:C	1:A:26:ILE:CD1	0.45	2.90	12	1
1:A:26:ILE:HD12	1:A:26:ILE:O	0.44	2.11	12	1
1:A:74:ASN:OD1	1:A:76:CYS:SG	0.44	2.74	13	1
1:A:111:ILE:O	1:A:111:ILE:CG2	0.43	2.65	7	1
1:A:95:PHE:HD1	1:A:99:CYS:HG	0.43	1.56	19	1
1:A:77:ASP:OD1	1:A:110:ARG:NH2	0.42	2.52	16	1
1:A:141:ILE:O	1:A:141:ILE:HG23	0.42	2.15	2	1
1:A:85:ARG:HD2	1:A:88:SER:O	0.41	2.15	9	1
1:A:76:CYS:HB2	1:A:79:LYS:O	0.41	2.15	5	1
1:A:114:GLU:O	1:A:118:CYS:SG	0.41	2.77	6	1
1:A:139:PHE:O	1:A:140:LYS:HB2	0.41	2.16	6	1
1:A:107:HIS:NE2	1:A:161:THR:OG1	0.41	2.50	4	1
1:A:141:ILE:O	1:A:141:ILE:CG2	0.41	2.64	20	1
1:A:178:TYR:CD1	1:A:178:TYR:C	0.40	2.99	10	1
1:A:149:LYS:NZ	1:A:181:LEU:OXT	0.40	2.47	15	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	148/170 (87%)	139±2 (94±1%)	8±2 (6±1%)	0±1 (0±1%)	37	78
All	All	2960/3400 (87%)	2785 (94%)	165 (6%)	10 (0%)	37	78

All 5 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	77	ASP	3
1	A	113	SER	3
1	A	165	ILE	2
1	A	146	ILE	1
1	A	39	GLY	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	139/160 (87%)	139±0 (100±0%)	0±0 (0±0%)	85	97
All	All	2780/3200 (87%)	2775 (100%)	5 (0%)	85	97

All 5 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	177	VAL	1
1	A	69	ILE	1
1	A	115	THR	1
1	A	26	ILE	1
1	A	129	TYR	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 47% for the well-defined parts and 47% 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	1071
Number of shifts mapped to atoms	1071
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	166	0.10 ± 0.15	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	161	0.83 ± 0.11	Should be checked
$^{13}\text{C}'$	146	0.02 ± 0.17	None needed (< 0.5 ppm)
^{15}N	151	0.91 ± 0.35	Should be applied

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 47%, i.e. 969 atoms were assigned a chemical shift out of a possible 2048. 0 out of 24 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	548/736 (74%)	136/295 (46%)	276/298 (93%)	136/143 (95%)
Sidechain	421/1140 (37%)	189/747 (25%)	232/357 (65%)	0/36 (0%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	0/172 (0%)	0/89 (0%)	0/82 (0%)	0/1 (0%)
Overall	969/2048 (47%)	325/1131 (29%)	508/737 (69%)	136/180 (76%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 47%, i.e. 1071 atoms were assigned a chemical shift out of a possible 2301. 0 out of 26 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	614/837 (73%)	151/335 (45%)	312/340 (92%)	151/162 (93%)
Sidechain	457/1251 (37%)	201/819 (25%)	256/395 (65%)	0/37 (0%)
Aromatic	0/213 (0%)	0/111 (0%)	0/101 (0%)	0/1 (0%)
Overall	1071/2301 (47%)	352/1265 (28%)	568/836 (68%)	151/200 (76%)

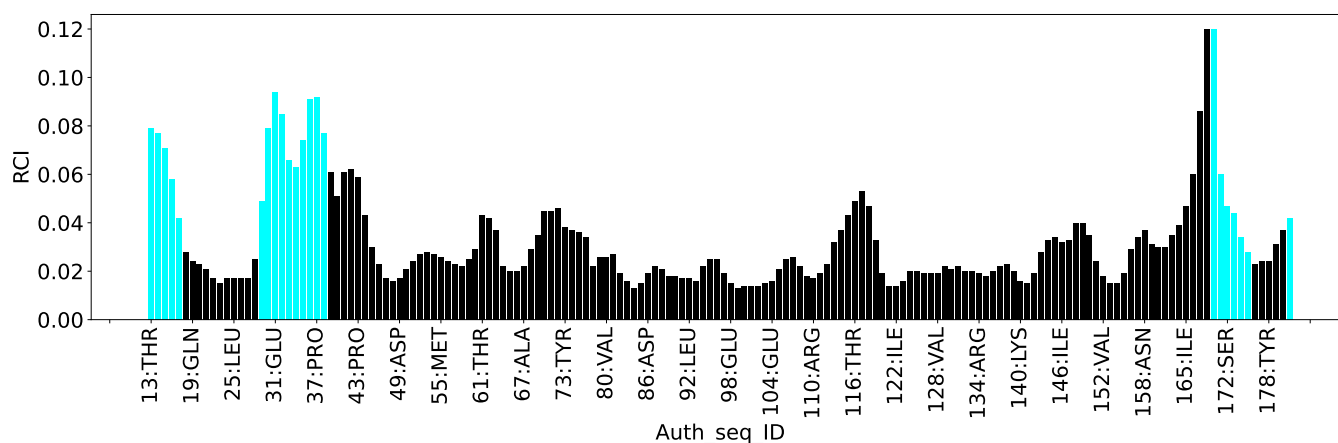
7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

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

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

Random coil index (RCI) for chain A:



8 NMR restraints analysis

8.1 Conformationally restricting restraints

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

Description	Value
Total distance restraints	538
Intra-residue ($ i-j =0$)	0
Sequential ($ i-j =1$)	109
Medium range ($ i-j >1$ and $ i-j <5$)	67
Long range ($ i-j \geq 5$)	362
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	3.2
Number of long range restraints per residue ¹	2.1

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

8.2 Residual restraint violations

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

8.2.1 Average number of distance violations per model

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

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	14.3	0.2
0.2-0.5 (Medium)	36.9	0.5
>0.5 (Large)	88.5	3.19

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

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

9 Distance violation analysis ⓘ

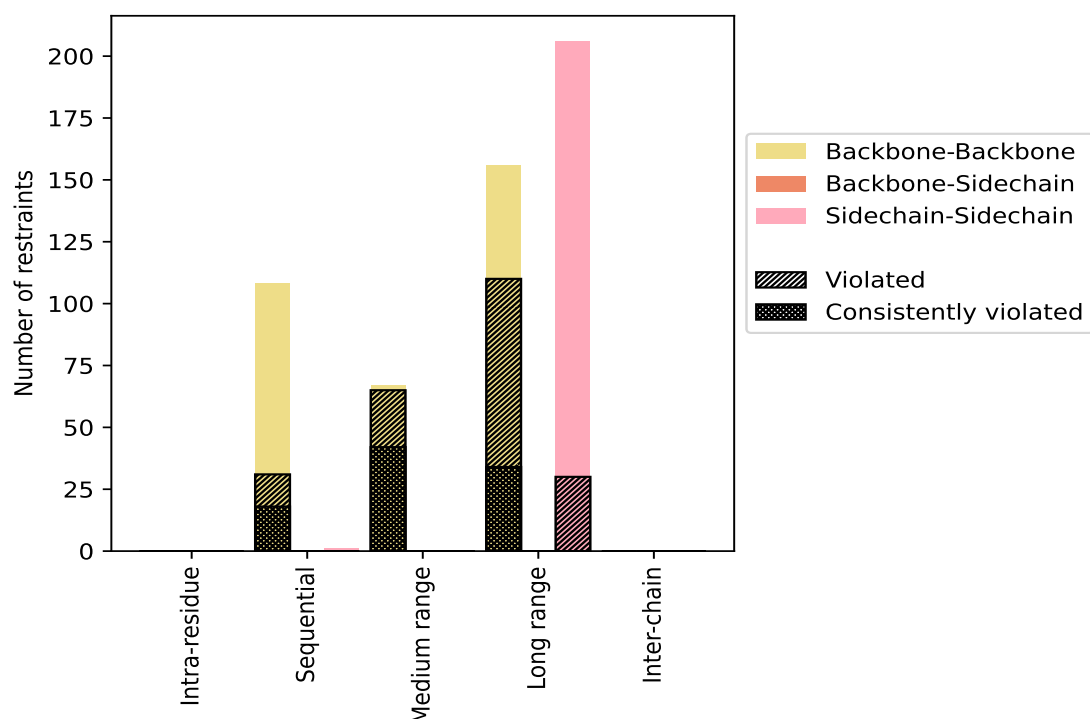
9.1 Summary of distance violations ⓘ

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

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($ i-j =0$)	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sequential ($ i-j =1$)	109	20.3	31	28.4	5.8	18	16.5	3.3
Backbone-Backbone	108	20.1	31	28.7	5.8	18	16.7	3.3
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	1	0.2	0	0.0	0.0	0	0.0	0.0
Medium range ($ i-j >1$ & $ i-j <5$)	67	12.5	65	97.0	12.1	42	62.7	7.8
Backbone-Backbone	67	12.5	65	97.0	12.1	42	62.7	7.8
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Long range ($ i-j \geq 5$)	362	67.3	140	38.7	26.0	34	9.4	6.3
Backbone-Backbone	156	29.0	110	70.5	20.4	34	21.8	6.3
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	206	38.3	30	14.6	5.6	0	0.0	0.0
Inter-chain	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Hydrogen bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	538	100.0	236	43.9	43.9	94	17.5	17.5
Backbone-Backbone	331	61.5	206	62.2	38.3	94	28.4	17.5
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	207	38.5	30	14.5	5.6	0	0.0	0.0

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

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



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

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

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

Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	0	22	52	71	0	145	0.7	2.53	0.45	0.68
2	0	21	55	61	0	137	0.69	2.82	0.43	0.6
3	0	22	58	61	0	141	0.67	2.35	0.42	0.63
4	0	21	57	71	0	149	0.66	2.31	0.4	0.63
5	0	22	56	62	0	140	0.67	2.23	0.42	0.59
6	0	25	53	57	0	135	0.72	2.57	0.47	0.61
7	0	20	53	55	0	128	0.75	2.71	0.47	0.7
8	0	24	53	64	0	141	0.74	3.19	0.52	0.71
9	0	22	55	62	0	139	0.72	2.37	0.46	0.64
10	0	21	51	64	0	136	0.68	2.65	0.41	0.62

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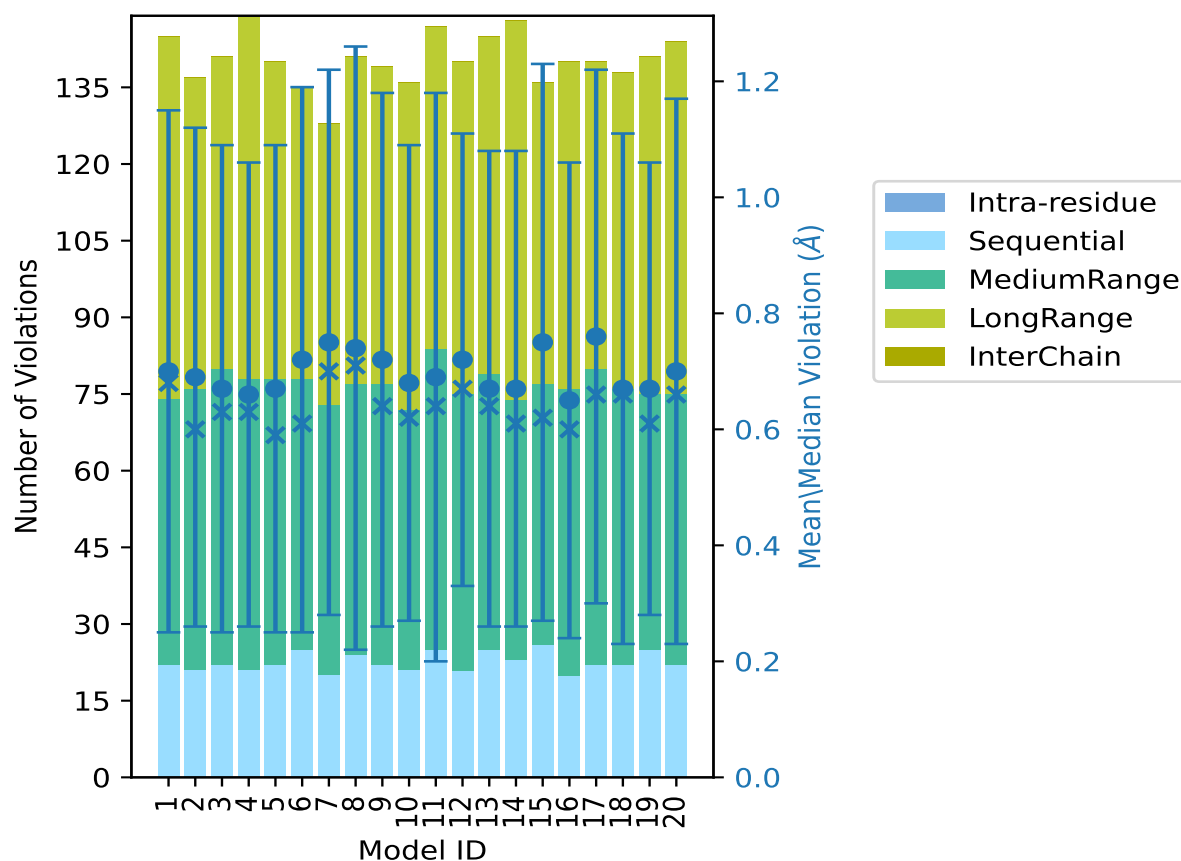
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
11	0	25	59	63	0	147	0.69	2.87	0.49	0.64
12	0	21	54	65	0	140	0.72	2.32	0.39	0.67
13	0	25	54	66	0	145	0.67	2.24	0.41	0.64
14	0	23	51	74	0	148	0.67	2.46	0.41	0.61
15	0	26	51	59	0	136	0.75	2.83	0.48	0.62
16	0	20	56	64	0	140	0.65	2.26	0.41	0.6
17	0	22	58	60	0	140	0.76	2.59	0.46	0.66
18	0	22	54	62	0	138	0.67	2.45	0.44	0.66
19	0	25	50	66	0	141	0.67	2.49	0.39	0.61
20	0	22	53	69	0	144	0.7	2.62	0.47	0.66

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

⁵Inter-chain restraints, ⁶Standard deviation

9.2.1 Bar graph : Distance Violation statistics for each model ⓘ



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

9.3 Distance violation statistics for the ensemble

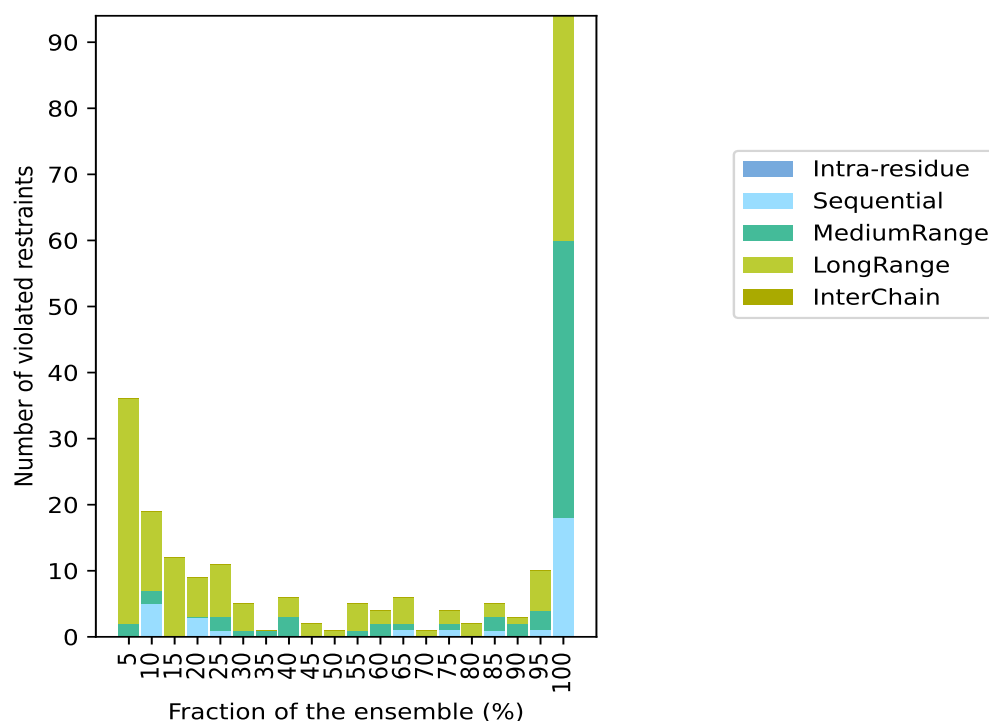
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 302(IR:0, SQ:78, MR:2, LR:222, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
0	0	2	34	0	36	1	5.0
0	5	2	12	0	19	2	10.0
0	0	0	12	0	12	3	15.0
0	3	0	6	0	9	4	20.0
0	1	2	8	0	11	5	25.0
0	0	1	4	0	5	6	30.0
0	0	1	0	0	1	7	35.0
0	0	3	3	0	6	8	40.0
0	0	0	2	0	2	9	45.0
0	0	0	1	0	1	10	50.0
0	0	1	4	0	5	11	55.0
0	0	2	2	0	4	12	60.0
0	1	1	4	0	6	13	65.0
0	0	0	1	0	1	14	70.0
0	1	1	2	0	4	15	75.0
0	0	0	2	0	2	16	80.0
0	1	2	2	0	5	17	85.0
0	0	2	1	0	3	18	90.0
0	1	3	6	0	10	19	95.0
0	18	42	34	0	94	20	100.0

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

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

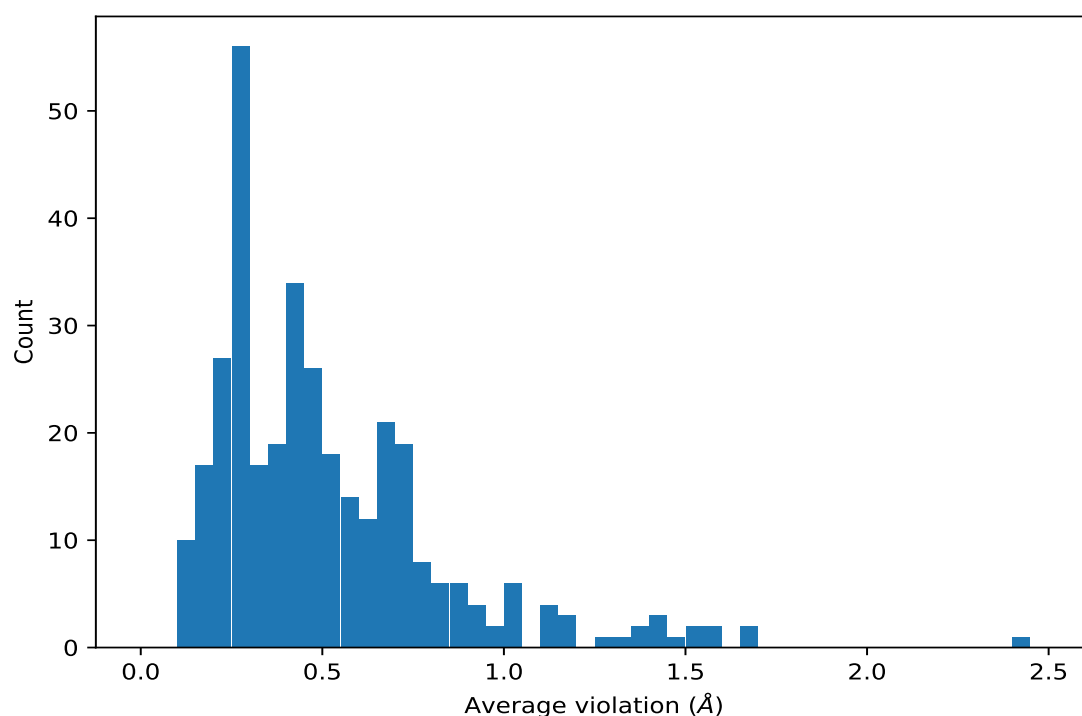
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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

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

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



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

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

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	20	2.42	0.2	2.38
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	20	1.7	0.42	1.69
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	20	1.66	0.64	1.28
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	20	1.55	0.42	1.54
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	20	1.55	0.42	1.54
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	20	1.51	0.34	1.54
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	20	1.5	0.28	1.5
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	20	1.45	0.04	1.45
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	20	1.35	0.51	1.41
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	20	1.3	0.17	1.33
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	20	1.29	0.1	1.32
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	20	1.2	0.51	1.26
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	20	1.2	0.51	1.26
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	20	1.19	0.26	1.13
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	20	1.15	0.17	1.18
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	20	1.15	0.17	1.18

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	20	1.11	0.41	1.0
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	20	1.1	0.04	1.1
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	20	1.04	0.26	0.98
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	20	1.04	0.26	0.98
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	20	1.04	0.64	0.66
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	20	1.03	0.1	1.05
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	20	1.0	0.16	1.02
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	20	1.0	0.1	1.02
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	20	0.96	0.11	0.94
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	20	0.96	0.31	0.9
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	20	0.95	0.1	0.94
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	20	0.95	0.11	0.96
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	20	0.92	0.1	0.94
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	20	0.92	0.1	0.9
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	20	0.89	0.1	0.88
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	20	0.88	0.04	0.88
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	20	0.88	0.1	0.89
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	20	0.88	0.1	0.89
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	20	0.87	0.31	0.82
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	20	0.87	0.08	0.87
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	20	0.84	0.09	0.84
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	20	0.83	0.21	0.92
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	20	0.83	0.09	0.82
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	20	0.82	0.14	0.76
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	20	0.81	0.06	0.85
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	20	0.8	0.1	0.82
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	20	0.79	0.01	0.79
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	20	0.79	0.13	0.8
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	20	0.78	0.16	0.8
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.76	0.16	0.78
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.76	0.07	0.76
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	20	0.76	0.11	0.75
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	20	0.76	0.04	0.78
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.73	0.07	0.74
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	20	0.7	0.04	0.7
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	20	0.69	0.14	0.74
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	20	0.69	0.14	0.66
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	20	0.68	0.21	0.77
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	20	0.68	0.21	0.77
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	20	0.68	0.04	0.69
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	20	0.67	0.14	0.6
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	20	0.67	0.14	0.6

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	20	0.65	0.14	0.67
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	20	0.64	0.13	0.64
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	20	0.64	0.13	0.64
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	20	0.62	0.07	0.62
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	20	0.62	0.09	0.61
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.61	0.16	0.62
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.61	0.16	0.62
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.61	0.16	0.62
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.61	0.07	0.62
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.61	0.07	0.62
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	20	0.6	0.12	0.6
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	20	0.6	0.07	0.59
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	20	0.58	0.11	0.6
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.58	0.07	0.58
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.58	0.07	0.58
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	20	0.57	0.01	0.58
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	20	0.57	0.06	0.57
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	20	0.56	0.13	0.57
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	20	0.56	0.08	0.57
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	20	0.48	0.11	0.5
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.46	0.16	0.48
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.46	0.16	0.48
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	20	0.43	0.11	0.44
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	20	0.43	0.11	0.44
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	20	0.42	0.06	0.42
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	20	0.42	0.06	0.42
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	20	0.42	0.21	0.32
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	20	0.41	0.08	0.42
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	20	0.38	0.13	0.36
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	20	0.38	0.06	0.4
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	20	0.34	0.01	0.34
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	20	0.33	0.03	0.32
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	20	0.3	0.03	0.3
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	20	0.23	0.06	0.24
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	20	0.21	0.03	0.21
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	20	0.18	0.03	0.18
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	19	1.43	0.17	1.39
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	19	1.43	0.17	1.39
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	19	0.64	0.45	0.5
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	19	0.6	0.02	0.6
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	19	0.55	0.25	0.46
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	19	0.43	0.15	0.42

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	19	0.42	0.2	0.39
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	19	0.4	0.25	0.31
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	19	0.4	0.25	0.31
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	19	0.37	0.15	0.38
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	18	0.42	0.16	0.42
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	18	0.42	0.26	0.34
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	18	0.3	0.11	0.3
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	17	0.6	0.36	0.46
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	17	0.58	0.2	0.69
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	17	0.4	0.26	0.31
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	17	0.37	0.14	0.36
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	17	0.3	0.1	0.29
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	16	0.31	0.13	0.28
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	16	0.31	0.13	0.28
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	15	1.49	0.49	1.69
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	15	0.45	0.26	0.33
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	15	0.31	0.12	0.29
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	15	0.2	0.07	0.18
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	14	0.75	0.38	0.75
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	13	1.35	0.04	1.36
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	13	0.37	0.12	0.37
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	13	0.36	0.25	0.31
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	13	0.36	0.25	0.31
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	13	0.28	0.1	0.25
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	13	0.28	0.1	0.25
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	12	0.69	0.34	0.64
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	12	0.69	0.34	0.64
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	12	0.48	0.09	0.46
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	12	0.43	0.22	0.4
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	11	0.37	0.25	0.3
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	11	0.37	0.25	0.3
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	11	0.24	0.08	0.21
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	11	0.21	0.06	0.2
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	11	0.18	0.08	0.19
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	10	0.5	0.43	0.26
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	10	0.5	0.43	0.26
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	10	0.5	0.43	0.26
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	10	0.5	0.43	0.26
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	10	0.5	0.43	0.26
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	10	0.5	0.43	0.26
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	10	0.5	0.43	0.26
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	10	0.5	0.43	0.26

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	10	0.5	0.43	0.26
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	9	0.23	0.08	0.23
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	9	0.23	0.08	0.23
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	8	0.62	0.17	0.68
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	8	0.59	0.35	0.52
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	8	0.47	0.17	0.53
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	8	0.47	0.17	0.53
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	8	0.14	0.01	0.14
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	8	0.14	0.01	0.14
(1,126)	1:86:A:ASP:H	1:88:A:SER:H	7	0.25	0.05	0.26
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	6	0.54	0.29	0.42
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	6	0.54	0.29	0.42
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	6	0.54	0.29	0.42
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	6	0.54	0.29	0.42
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	6	0.54	0.29	0.42
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	6	0.54	0.29	0.42
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	6	0.54	0.29	0.42
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	6	0.54	0.29	0.42
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	6	0.54	0.29	0.42
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	6	0.48	0.29	0.36
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	6	0.48	0.29	0.36
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	6	0.48	0.29	0.36
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	6	0.48	0.29	0.36
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	6	0.48	0.29	0.36
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	6	0.48	0.29	0.36
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	6	0.48	0.29	0.36
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	6	0.48	0.29	0.36
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	6	0.48	0.29	0.36
(1,3)	1:119:A:LEU:H	1:163:A:PHE:H	6	0.33	0.12	0.34
(1,104)	1:84:A:ILE:H	1:91:A:ILE:H	6	0.17	0.04	0.18
(1,159)	1:46:A:THR:H	1:49:A:ASP:H	6	0.14	0.04	0.13
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	5	0.69	0.38	0.85
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	5	0.69	0.38	0.85
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	5	0.69	0.38	0.85
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	5	0.69	0.38	0.85
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	5	0.69	0.38	0.85
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	5	0.69	0.38	0.85
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	5	0.69	0.38	0.85
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	5	0.69	0.38	0.85
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	5	0.69	0.38	0.85
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	5	0.44	0.28	0.34
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	5	0.44	0.28	0.34

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	5	0.44	0.28	0.34
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	5	0.44	0.28	0.34
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	5	0.44	0.28	0.34
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	5	0.44	0.28	0.34
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	5	0.44	0.28	0.34
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	5	0.44	0.28	0.34
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	5	0.44	0.28	0.34
(1,40)	1:124:A:ASP:H	1:125:A:ALA:H	5	0.36	0.46	0.12
(1,178)	1:90:A:GLN:H	1:177:A:VAL:H	5	0.26	0.11	0.34
(1,107)	1:67:A:ALA:H	1:120:A:MET:H	5	0.23	0.11	0.23
(3,2)	1:119:A:LEU:H	1:163:A:PHE:H	5	0.23	0.07	0.21
(3,120)	1:119:A:LEU:H	1:163:A:PHE:H	5	0.23	0.07	0.21
(3,16)	1:131:A:THR:H	1:135:A:ALA:H	5	0.17	0.06	0.14
(3,134)	1:131:A:THR:H	1:135:A:ALA:H	5	0.17	0.06	0.14
(3,8)	1:126:A:ASP:H	1:156:A:ASN:H	5	0.14	0.05	0.12
(3,126)	1:126:A:ASP:H	1:156:A:ASN:H	5	0.14	0.05	0.12
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	4	0.72	0.29	0.84
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	4	0.72	0.29	0.84
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	4	0.72	0.29	0.84
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	4	0.72	0.29	0.84
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	4	0.72	0.29	0.84
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	4	0.72	0.29	0.84
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	4	0.72	0.29	0.84
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	4	0.72	0.29	0.84
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	4	0.72	0.29	0.84
(1,155)	1:30:A:SER:H	1:31:A:GLU:H	4	0.7	0.4	0.57
(1,53)	1:30:A:SER:H	1:31:A:GLU:H	4	0.64	0.4	0.5
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	4	0.46	0.26	0.48
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	4	0.46	0.26	0.48
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	4	0.46	0.26	0.48
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	4	0.46	0.26	0.48
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	4	0.46	0.26	0.48
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	4	0.46	0.26	0.48
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	4	0.46	0.26	0.48
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	4	0.46	0.26	0.48
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	4	0.46	0.26	0.48
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	4	0.35	0.27	0.26
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	4	0.35	0.27	0.26
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	4	0.35	0.27	0.26
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	4	0.35	0.27	0.26
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	4	0.35	0.27	0.26
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	4	0.35	0.27	0.26

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	4	0.35	0.27	0.26
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	4	0.35	0.27	0.26
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	4	0.35	0.27	0.26
(1,110)	1:45:A:LEU:H	1:124:A:ASP:H	4	0.22	0.08	0.2
(1,182)	1:108:A:ASN:H	1:109:A:CYS:H	4	0.14	0.02	0.14
(3,42)	1:84:A:ILE:H	1:91:A:ILE:H	4	0.14	0.02	0.14
(3,160)	1:84:A:ILE:H	1:91:A:ILE:H	4	0.14	0.02	0.14
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	3	0.74	0.1	0.78
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	3	0.74	0.1	0.78
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	3	0.74	0.1	0.78
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	3	0.74	0.1	0.78
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	3	0.74	0.1	0.78
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	3	0.74	0.1	0.78
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	3	0.74	0.1	0.78
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	3	0.74	0.1	0.78
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	3	0.74	0.1	0.78
(1,209)	1:21:A:TRP:H	1:87:A:SER:H	3	0.47	0.12	0.39
(3,52)	1:21:A:TRP:H	1:87:A:SER:H	3	0.32	0.12	0.24
(3,170)	1:21:A:TRP:H	1:87:A:SER:H	3	0.32	0.12	0.24
(1,118)	1:68:A:LYS:H	1:85:A:ARG:H	3	0.3	0.2	0.19
(1,149)	1:68:A:LYS:H	1:85:A:ARG:H	3	0.3	0.2	0.19
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD11	3	0.27	0.1	0.27
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD12	3	0.27	0.1	0.27
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD13	3	0.27	0.1	0.27
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD11	3	0.27	0.1	0.27
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD12	3	0.27	0.1	0.27
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD13	3	0.27	0.1	0.27
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD11	3	0.27	0.1	0.27
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD12	3	0.27	0.1	0.27
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD13	3	0.27	0.1	0.27
(1,94)	1:121:A:ILE:H	1:163:A:PHE:H	3	0.22	0.05	0.23
(3,46)	1:90:A:GLN:H	1:177:A:VAL:H	3	0.2	0.01	0.2
(3,164)	1:90:A:GLN:H	1:177:A:VAL:H	3	0.2	0.01	0.2
(1,75)	1:90:A:GLN:H	1:177:A:VAL:H	3	0.13	0.01	0.13
(1,33)	1:94:A:GLU:H	1:155:A:LEU:H	3	0.13	0.0	0.13
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	2	0.42	0.23	0.42
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	2	0.42	0.23	0.42
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	2	0.42	0.23	0.42
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	2	0.42	0.23	0.42
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	2	0.42	0.23	0.42
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	2	0.42	0.23	0.42
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	2	0.42	0.23	0.42

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	2	0.42	0.23	0.42
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	2	0.42	0.23	0.42
(1,15)	1:180:A:TYR:H	1:181:A:LEU:H	2	0.4	0.2	0.4
(1,129)	1:180:A:TYR:H	1:181:A:LEU:H	2	0.4	0.2	0.4
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	2	0.36	0.23	0.36
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	2	0.36	0.23	0.36
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	2	0.36	0.23	0.36
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	2	0.36	0.23	0.36
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	2	0.36	0.23	0.36
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	2	0.36	0.23	0.36
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	2	0.36	0.23	0.36
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	2	0.36	0.23	0.36
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	2	0.36	0.23	0.36
(1,83)	1:138:A:HIS:H	1:140:A:LYS:H	2	0.28	0.07	0.28
(1,185)	1:138:A:HIS:H	1:140:A:LYS:H	2	0.28	0.07	0.28
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG21	2	0.27	0.05	0.27
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG22	2	0.27	0.05	0.27
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG23	2	0.27	0.05	0.27
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG21	2	0.27	0.05	0.27
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG22	2	0.27	0.05	0.27
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG23	2	0.27	0.05	0.27
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG21	2	0.27	0.05	0.27
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG22	2	0.27	0.05	0.27
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG23	2	0.27	0.05	0.27
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG11	2	0.26	0.08	0.26
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG12	2	0.26	0.08	0.26
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG13	2	0.26	0.08	0.26
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG11	2	0.26	0.08	0.26
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG12	2	0.26	0.08	0.26
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG13	2	0.26	0.08	0.26
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG11	2	0.26	0.08	0.26
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG12	2	0.26	0.08	0.26
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG13	2	0.26	0.08	0.26
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD11	2	0.26	0.11	0.26
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD12	2	0.26	0.11	0.26
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD13	2	0.26	0.11	0.26
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD11	2	0.26	0.11	0.26
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD12	2	0.26	0.11	0.26
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD13	2	0.26	0.11	0.26
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD11	2	0.26	0.11	0.26
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD12	2	0.26	0.11	0.26
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD13	2	0.26	0.11	0.26

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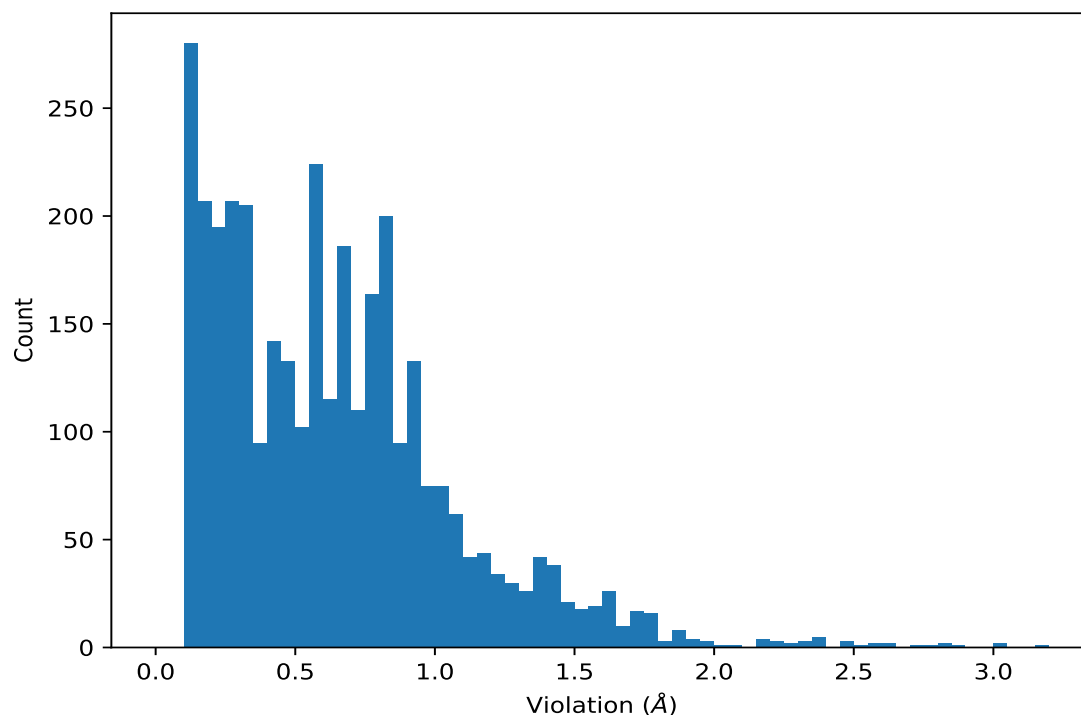
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG21	2	0.26	0.05	0.26
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG22	2	0.26	0.05	0.26
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG23	2	0.26	0.05	0.26
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG21	2	0.26	0.05	0.26
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG22	2	0.26	0.05	0.26
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG23	2	0.26	0.05	0.26
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG21	2	0.26	0.05	0.26
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG22	2	0.26	0.05	0.26
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG23	2	0.26	0.05	0.26
(1,41)	1:42:A:ILE:H	1:125:A:ALA:H	2	0.24	0.08	0.24
(1,88)	1:42:A:ILE:H	1:125:A:ALA:H	2	0.24	0.08	0.24
(1,125)	1:88:A:SER:H	1:89:A:HIS:H	2	0.23	0.01	0.23
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD11	2	0.22	0.06	0.22
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD12	2	0.22	0.06	0.22
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD13	2	0.22	0.06	0.22
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD11	2	0.22	0.06	0.22
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD12	2	0.22	0.06	0.22
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD13	2	0.22	0.06	0.22
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD11	2	0.22	0.06	0.22
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD12	2	0.22	0.06	0.22
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD13	2	0.22	0.06	0.22
(1,55)	1:88:A:SER:H	1:89:A:HIS:H	2	0.2	0.01	0.2
(3,32)	1:67:A:ALA:H	1:120:A:MET:H	2	0.19	0.07	0.19
(3,150)	1:67:A:ALA:H	1:120:A:MET:H	2	0.19	0.07	0.19
(1,203)	1:50:A:LEU:H	1:51:A:ALA:H	2	0.18	0.02	0.18
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD11	2	0.18	0.06	0.18
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD12	2	0.18	0.06	0.18
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD13	2	0.18	0.06	0.18
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD11	2	0.18	0.06	0.18
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD12	2	0.18	0.06	0.18
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD13	2	0.18	0.06	0.18
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD11	2	0.18	0.06	0.18
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD12	2	0.18	0.06	0.18
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD13	2	0.18	0.06	0.18

¹Number of violated models, ²Standard deviation

9.5 All violated distance restraints [i](#)

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

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



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

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

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	8	3.19
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	8	3.04
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	8	3.04
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	11	2.87
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	15	2.83
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	2	2.82
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	11	2.77
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	7	2.71
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	10	2.65
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	20	2.62

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	17	2.59
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	6	2.57
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	1	2.53
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	19	2.49
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	14	2.46
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	18	2.45
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	1	2.4
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	8	2.38
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	9	2.37
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	15	2.36
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	3	2.35
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	12	2.32
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	20	2.32
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	4	2.31
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	7	2.28
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	16	2.26
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	13	2.24
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	5	2.23
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	15	2.21
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	18	2.18
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	20	2.17
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	20	2.17
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	11	2.15
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	7	2.09
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	6	2.03
(1,215)	1:141:A:ILE:H	1:143:A:LEU:H	17	1.98
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	9	1.97
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	17	1.97
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	20	1.94
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	9	1.94
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	6	1.94
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	1	1.91
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	17	1.9
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	15	1.9
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	17	1.89
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	15	1.88
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	8	1.86
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	20	1.86
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	16	1.86
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	6	1.86
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	7	1.82
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	2	1.8

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	14	1.8
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	20	1.79
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	9	1.79
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	20	1.79
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	9	1.79
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	10	1.79
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	13	1.79
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	4	1.78
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	5	1.78
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	16	1.78
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	9	1.76
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	17	1.76
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	17	1.75
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	15	1.75
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	17	1.75
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	15	1.75
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	3	1.75
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	17	1.74
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	17	1.74
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	6	1.74
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	8	1.74
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	18	1.74
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	15	1.73
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	15	1.73
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	6	1.73
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	2	1.72
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	11	1.72
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	15	1.72
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	16	1.71
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	6	1.71
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	16	1.71
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	6	1.71
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	18	1.71
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	14	1.71
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	9	1.7
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	13	1.7
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	2	1.69
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	1	1.69
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	5	1.68
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	7	1.67
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	7	1.67
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	18	1.67

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	19	1.66
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	19	1.66
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	2	1.65
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	14	1.65
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	2	1.65
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	14	1.65
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	4	1.65
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	10	1.65
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	3	1.65
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	13	1.64
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	13	1.64
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	11	1.64
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	13	1.64
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	4	1.63
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	4	1.63
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	9	1.63
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	10	1.62
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	17	1.62
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	1	1.62
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	9	1.61
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	17	1.61
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	9	1.61
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	17	1.61
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	20	1.61
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	3	1.6
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	3	1.6
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	7	1.6
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	17	1.6
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	6	1.59
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	6	1.59
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	11	1.59
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	12	1.59
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	12	1.59
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	2	1.57
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	11	1.57
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	2	1.57
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	11	1.57
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	7	1.57
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	20	1.57
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	18	1.56
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	18	1.56
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	3	1.56

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	19	1.56
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	1	1.56
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	5	1.56
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	5	1.55
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	7	1.55
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	13	1.54
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	4	1.54
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	14	1.54
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	6	1.53
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	17	1.53
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	18	1.52
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	18	1.52
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	11	1.52
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	12	1.52
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	12	1.52
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	19	1.51
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	19	1.51
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	4	1.5
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	10	1.5
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	4	1.5
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	10	1.5
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	11	1.5
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	12	1.5
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	7	1.49
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	5	1.49
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	9	1.49
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	13	1.49
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	6	1.49
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	9	1.48
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	6	1.48
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	16	1.48
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	18	1.48
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	10	1.47
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	1	1.47
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	10	1.47
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	1	1.47
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	2	1.47
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	15	1.47
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	4	1.47
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	19	1.47
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	11	1.46
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	14	1.46

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	15	1.46
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	5	1.46
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	7	1.45
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	7	1.45
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	4	1.45
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	3	1.45
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	2	1.45
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	10	1.45
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	11	1.44
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	12	1.44
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	11	1.44
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	12	1.44
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	17	1.44
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	4	1.44
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	1	1.44
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	3	1.44
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	7	1.44
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	14	1.44
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	20	1.44
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	8	1.44
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	13	1.43
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	1	1.43
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	16	1.43
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	18	1.43
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	15	1.43
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	8	1.43
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	17	1.43
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	7	1.42
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	20	1.42
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	7	1.42
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	20	1.42
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	5	1.42
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	19	1.41
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	1	1.41
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	5	1.41
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	19	1.41
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	1	1.41
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	5	1.41
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	13	1.41
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	14	1.41
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	7	1.4
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	7	1.4

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	3	1.4
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	5	1.4
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	12	1.4
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	12	1.4
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	8	1.4
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	18	1.4
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	1	1.4
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	9	1.4
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	19	1.4
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	4	1.39
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	4	1.39
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	15	1.39
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	16	1.39
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	5	1.39
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	6	1.39
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	11	1.39
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	6	1.38
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	6	1.38
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	9	1.38
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	8	1.38
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	15	1.38
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	12	1.37
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	12	1.37
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	12	1.37
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	12	1.37
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	14	1.37
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	8	1.37
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	6	1.36
(1,155)	1:30:A:SER:H	1:31:A:GLU:H	8	1.36
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	3	1.36
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	9	1.36
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	13	1.36
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	12	1.36
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	16	1.35
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	15	1.35
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	18	1.35
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	11	1.35
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	1	1.35
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	9	1.35
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	20	1.35
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	5	1.34
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	5	1.34

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	1	1.34
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	2	1.34
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	15	1.34
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	9	1.33
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	9	1.33
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	2	1.32
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	15	1.32
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	2	1.32
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	15	1.32
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	19	1.32
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	7	1.32
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	3	1.32
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	11	1.31
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	11	1.31
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	14	1.31
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	6	1.31
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	12	1.31
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	3	1.31
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	17	1.31
(1,43)	1:167:A:GLN:H	1:168:A:VAL:H	11	1.31
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	3	1.3
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	3	1.3
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	11	1.3
(1,53)	1:30:A:SER:H	1:31:A:GLU:H	8	1.3
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	17	1.29
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	17	1.29
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	13	1.28
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	1	1.28
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	16	1.28
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	18	1.28
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	13	1.28
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	1	1.28
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	16	1.28
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	18	1.28
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	17	1.28
(1,40)	1:124:A:ASP:H	1:125:A:ALA:H	8	1.28
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	4	1.27
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	19	1.27
(1,73)	1:35:A:TYR:H	1:36:A:HIS:H	20	1.27
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	12	1.27
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	3	1.27
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	14	1.27

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	13	1.26
(3,153)	1:124:A:ASP:H	1:159:A:GLN:H	14	1.26
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	13	1.26
(3,35)	1:124:A:ASP:H	1:159:A:GLN:H	14	1.26
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	3	1.25
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	5	1.25
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	12	1.25
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	3	1.25
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	5	1.25
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	12	1.25
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	8	1.25
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	13	1.25
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	15	1.24
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	5	1.24
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	15	1.24
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	5	1.24
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	1	1.24
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	3	1.24
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	20	1.24
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	9	1.23
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	8	1.23
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	9	1.23
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	8	1.23
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	14	1.23
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	14	1.23
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	14	1.23
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	14	1.23
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	14	1.23
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	14	1.23
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	14	1.23
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	14	1.23
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	14	1.23
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	18	1.23
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	19	1.23
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	10	1.23
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	16	1.23
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	10	1.23
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	2	1.22
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	10	1.22
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	12	1.22
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	13	1.22
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	14	1.22

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	16	1.22
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	3	1.21
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	3	1.21
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	5	1.21
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	16	1.2
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	11	1.2
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	16	1.2
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	11	1.2
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	16	1.2
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	7	1.2
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	5	1.2
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	1	1.19
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	1	1.19
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	3	1.19
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	10	1.19
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	4	1.19
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	19	1.19
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	6	1.18
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	6	1.18
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	6	1.18
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	6	1.18
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	6	1.18
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	6	1.18
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	6	1.18
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	6	1.18
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	6	1.18
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	2	1.18
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	10	1.18
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	5	1.18
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	6	1.18
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	10	1.17
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	13	1.17
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	13	1.17
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	14	1.16
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	6	1.16
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	14	1.16
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	6	1.16
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	2	1.16
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	14	1.16
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	5	1.16
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	16	1.16
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	19	1.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	3	1.16
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	10	1.16
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	8	1.16
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	9	1.16
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	17	1.16
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	9	1.16
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	14	1.15
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	19	1.15
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	12	1.15
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	15	1.15
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	1	1.14
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	3	1.14
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	9	1.14
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	13	1.14
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	9	1.13
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	8	1.13
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	19	1.13
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	2	1.13
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	13	1.13
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	6	1.13
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	16	1.13
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	18	1.13
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	4	1.12
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	19	1.12
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	12	1.12
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	4	1.12
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	19	1.12
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	12	1.12
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	7	1.12
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	8	1.12
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	11	1.12
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	4	1.12
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	19	1.12
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	11	1.11
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	3	1.11
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	15	1.11
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	9	1.11
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	8	1.1
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	13	1.1
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	8	1.1
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	13	1.1
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	20	1.1

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	9	1.1
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	13	1.1
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	2	1.1
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	5	1.1
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	2	1.1
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	10	1.1
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	1	1.09
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	1	1.09
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	2	1.09
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	6	1.09
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	8	1.09
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	7	1.09
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	11	1.09
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	4	1.09
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	16	1.09
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	5	1.09
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	1	1.09
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	3	1.09
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	7	1.09
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	14	1.09
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	20	1.09
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	16	1.09
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	6	1.09
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	18	1.08
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	10	1.08
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	14	1.08
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	14	1.08
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	14	1.08
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	14	1.08
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	14	1.08
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	14	1.08
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	14	1.08
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	14	1.08
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	14	1.08
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	18	1.08
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	10	1.08
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	12	1.08
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	10	1.08
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	11	1.08
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	10	1.08
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	9	1.08
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	10	1.08

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	15	1.08
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	8	1.08
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	17	1.08
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	9	1.08
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	18	1.08
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	19	1.07
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	20	1.07
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	5	1.07
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	5	1.07
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	15	1.07
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	14	1.06
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	12	1.06
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	17	1.06
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	5	1.06
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	7	1.05
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	5	1.05
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	7	1.05
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	5	1.05
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	1	1.05
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	17	1.05
(1,190)	1:87:A:SER:H	1:89:A:HIS:H	8	1.05
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	10	1.05
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	11	1.05
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	7	1.05
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	17	1.05
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	9	1.05
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	13	1.04
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	4	1.04
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	19	1.04
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	14	1.04
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	12	1.04
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	19	1.04
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	8	1.04
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	14	1.04
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	12	1.04
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	17	1.04
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	3	1.04
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	7	1.04
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	20	1.04
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	20	1.04
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	2	1.03
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	10	1.03

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	6	1.03
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	6	1.03
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	6	1.03
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	6	1.03
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	6	1.03
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	6	1.03
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	6	1.03
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	6	1.03
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	6	1.03
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	2	1.03
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	10	1.03
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	6	1.03
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	11	1.03
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	15	1.03
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	14	1.03
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	19	1.03
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	20	1.03
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	20	1.03
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	10	1.02
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	10	1.02
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	9	1.02
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	13	1.02
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	1	1.02
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	17	1.02
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	9	1.02
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	15	1.02
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	6	1.02
(1,52)	1:99:A:CYS:H	1:101:A:SER:H	18	1.02
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	11	1.02
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	17	1.02
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	19	1.01
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	8	1.01
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	9	1.01
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	19	1.01
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	8	1.01
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	9	1.01
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	5	1.01
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	1	1.01
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	4	1.01
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	14	1.01
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	3	1.01
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	16	1.01

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	11	1.01
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	7	1.01
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	14	1.0
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	14	1.0
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	13	1.0
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	13	1.0
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	13	1.0
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	13	1.0
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	13	1.0
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	13	1.0
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	13	1.0
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	13	1.0
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	13	1.0
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	15	1.0
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	2	1.0
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	10	1.0
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	17	1.0
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	3	0.99
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	3	0.99
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	10	0.99
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	4	0.99
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	2	0.99
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	19	0.99
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	1	0.99
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	20	0.99
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	6	0.99
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	16	0.99
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	9	0.99
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	8	0.98
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	13	0.98
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	8	0.98
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	13	0.98
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	14	0.98
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	14	0.98
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	14	0.98
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	14	0.98
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	14	0.98
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	14	0.98
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	14	0.98
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	14	0.98
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	14	0.98
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	14	0.98

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	14	0.98
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	14	0.98
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	14	0.98
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	14	0.98
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	14	0.98
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	14	0.98
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	14	0.98
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	14	0.98
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	14	0.98
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	3	0.98
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	15	0.98
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	17	0.98
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	3	0.98
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	20	0.98
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	20	0.98
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	13	0.98
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	17	0.98
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	3	0.98
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	17	0.98
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	10	0.98
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	10	0.98
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	11	0.98
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	5	0.97
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	12	0.97
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	2	0.97
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	8	0.97
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	18	0.97
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	11	0.97
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	16	0.97
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	10	0.97
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	2	0.97
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	8	0.97
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	18	0.97
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	9	0.97
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	1	0.97
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	8	0.97
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	1	0.97
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	1	0.97
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	12	0.97
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	17	0.97
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	7	0.96
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	18	0.96

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	18	0.96
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	2	0.96
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	5	0.96
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	11	0.96
(1,54)	1:167:A:GLN:H	1:168:A:VAL:H	11	0.96
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	4	0.96
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	2	0.96
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	17	0.96
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	18	0.95
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	17	0.95
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	11	0.95
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	11	0.95
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	11	0.95
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	4	0.95
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	2	0.94
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	8	0.94
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	4	0.94
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	16	0.94
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	5	0.94
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	2	0.94
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	8	0.94
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	4	0.94
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	16	0.94
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	5	0.94
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	8	0.94
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	7	0.94
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	16	0.94
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	12	0.94
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	20	0.94
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	5	0.94
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	7	0.94
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	16	0.94
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	4	0.94
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	12	0.94
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	10	0.94
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	9	0.94
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	1	0.94
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	3	0.94
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	9	0.94
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	19	0.94
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	17	0.94
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	15	0.94

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	14	0.94
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	15	0.94
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	9	0.94
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	7	0.94
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	12	0.93
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	18	0.93
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	12	0.93
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	18	0.93
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	6	0.93
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	6	0.93
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	6	0.93
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	6	0.93
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	6	0.93
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	6	0.93
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	6	0.93
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	6	0.93
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	6	0.93
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	2	0.93
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	19	0.93
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	6	0.93
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	6	0.93
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	15	0.93
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	8	0.93
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	18	0.93
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	8	0.93
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	12	0.93
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	4	0.93
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	7	0.93
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	7	0.93
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	1	0.93
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	2	0.93
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	12	0.93
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	17	0.93
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	2	0.93
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	9	0.93
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	1	0.93
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	15	0.92
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	15	0.92
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	14	0.92
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	14	0.92
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	14	0.92
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	14	0.92

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	14	0.92
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	14	0.92
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	14	0.92
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	14	0.92
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	14	0.92
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	20	0.92
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	14	0.92
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	13	0.92
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	19	0.92
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	18	0.92
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	13	0.92
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	12	0.92
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	16	0.92
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	13	0.92
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	14	0.92
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	14	0.92
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	7	0.92
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	5	0.91
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	2	0.91
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	3	0.91
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	7	0.91
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	1	0.91
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	20	0.91
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	8	0.91
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	5	0.91
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	6	0.91
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	1	0.91
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	9	0.91
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	19	0.91
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	10	0.91
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	16	0.91
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	19	0.91
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	20	0.91
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	15	0.91
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	7	0.91
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	4	0.91
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	1	0.91
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	5	0.91
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	8	0.91
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	10	0.91
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	17	0.9
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	17	0.9

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	18	0.9
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	12	0.9
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	16	0.9
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	16	0.9
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	3	0.9
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	11	0.9
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	11	0.9
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	1	0.9
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	16	0.9
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	18	0.9
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	8	0.9
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	9	0.9
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	10	0.9
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	5	0.9
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	6	0.9
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	8	0.89
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	14	0.89
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	12	0.89
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.89
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	8	0.89
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	14	0.89
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	12	0.89
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.89
(1,212)	1:138:A:HIS:H	1:141:A:ILE:H	20	0.89
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	20	0.89
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	9	0.89
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.89
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	13	0.89
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	8	0.89
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	1	0.89
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	18	0.89
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	1	0.89
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	7	0.89
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	2	0.89
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	20	0.89
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	6	0.88
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	19	0.88
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	20	0.88
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	6	0.88
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	19	0.88
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	20	0.88
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	12	0.88

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	12	0.88
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	12	0.88
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	12	0.88
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	12	0.88
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	12	0.88
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	12	0.88
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	12	0.88
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	12	0.88
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	4	0.88
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	12	0.88
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	8	0.88
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	3	0.88
(1,90)	1:42:A:ILE:H	1:124:A:ASP:H	20	0.88
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	2	0.88
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	11	0.88
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	18	0.88
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	18	0.88
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	11	0.88
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	15	0.88
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	19	0.88
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	15	0.87
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	6	0.87
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	6	0.87
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	15	0.87
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	20	0.87
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	14	0.87
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	12	0.87
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	16	0.87
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	15	0.87
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	20	0.87
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	12	0.87
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	5	0.87
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	6	0.87
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	9	0.87
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	16	0.87
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	13	0.87
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	4	0.87
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	8	0.87
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	18	0.87
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	18	0.87
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	7	0.86
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	7	0.86

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	12	0.86
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	16	0.86
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	1	0.86
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	12	0.86
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	13	0.86
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	15	0.86
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	16	0.86
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	1	0.86
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	12	0.86
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	20	0.86
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	15	0.86
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	3	0.86
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	7	0.86
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	10	0.86
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	4	0.86
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	5	0.86
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	6	0.86
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	17	0.86
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	2	0.86
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	4	0.86
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	11	0.86
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	1	0.86
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	2	0.86
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	8	0.86
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	4	0.86
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	19	0.86
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	13	0.85
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	13	0.85
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	13	0.85
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	13	0.85
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	13	0.85
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	13	0.85
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	13	0.85
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	13	0.85
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	13	0.85
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	6	0.85
(1,202)	1:50:A:LEU:H	1:52:A:LEU:H	13	0.85
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	1	0.85
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	2	0.85
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	5	0.85
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	7	0.85
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	9	0.85

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	11	0.85
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	12	0.85
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	19	0.85
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	1	0.85
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	2	0.85
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	15	0.85
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	1	0.85
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	7	0.85
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	16	0.85
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	18	0.85
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	10	0.85
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	16	0.85
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	5	0.85
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	9	0.85
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	12	0.85
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	13	0.85
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	18	0.85
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	14	0.85
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	16	0.85
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	6	0.84
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	16	0.84
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	9	0.84
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	6	0.84
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	16	0.84
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	9	0.84
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	3	0.84
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	10	0.84
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	16	0.84
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	2	0.84
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	10	0.84
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	6	0.84
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	8	0.84
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	20	0.84
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	2	0.84
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	17	0.84
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	14	0.84
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	10	0.84
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	16	0.84
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	12	0.84
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	15	0.84
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	3	0.84
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	13	0.84

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	19	0.84
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	8	0.84
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	11	0.83
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	14	0.83
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	14	0.83
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	14	0.83
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	14	0.83
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	14	0.83
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	14	0.83
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	14	0.83
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	14	0.83
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	14	0.83
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	14	0.83
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	14	0.83
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	14	0.83
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	14	0.83
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	14	0.83
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	14	0.83
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	14	0.83
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	14	0.83
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	14	0.83
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	11	0.83
(1,214)	1:137:A:SER:H	1:141:A:ILE:H	20	0.83
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	5	0.83
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	12	0.83
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	17	0.83
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	4	0.83
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	14	0.83
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	6	0.83
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	9	0.83
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	7	0.83
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	5	0.83
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	19	0.83
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	6	0.83
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	20	0.83
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	13	0.83
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	13	0.83
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	4	0.83
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	8	0.83
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	4	0.83
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	11	0.83
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	1	0.83

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	8	0.82
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	1	0.82
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	8	0.82
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	1	0.82
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	12	0.82
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	12	0.82
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	12	0.82
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	12	0.82
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	12	0.82
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	12	0.82
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	12	0.82
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	12	0.82
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	12	0.82
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	4	0.82
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	8	0.82
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	1	0.82
(1,140)	1:50:A:LEU:H	1:52:A:LEU:H	13	0.82
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	8	0.82
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	17	0.82
(1,102)	1:26:A:ILE:H	1:28:A:ASP:H	16	0.82
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	4	0.82
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	7	0.82
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	15	0.82
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	13	0.82
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	13	0.82
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	7	0.82
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	14	0.82
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	18	0.82
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	20	0.82
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	10	0.82
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	16	0.82
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	12	0.82
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	13	0.82
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	18	0.82
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	2	0.81
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	17	0.81
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	2	0.81
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	17	0.81
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	10	0.81
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	18	0.81
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	20	0.81
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	17	0.81

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	2	0.81
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	2	0.81
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	7	0.81
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	17	0.81
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	12	0.81
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	1	0.81
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	8	0.81
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	12	0.81
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	12	0.81
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	19	0.81
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	3	0.81
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	15	0.81
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	19	0.81
(1,25)	1:103:A:PHE:H	1:106:A:THR:H	20	0.81
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	13	0.81
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	11	0.81
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	17	0.8
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	11	0.8
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	4	0.8
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	17	0.8
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	11	0.8
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	4	0.8
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	16	0.8
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	16	0.8
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	16	0.8
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	16	0.8
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	16	0.8
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	16	0.8
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	16	0.8
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	16	0.8
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	16	0.8
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	4	0.8
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	17	0.8
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	2	0.8
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	2	0.8
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	6	0.8
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	7	0.8
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	4	0.8
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	17	0.8
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	4	0.8
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	7	0.8
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	1	0.8

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	3	0.8
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	8	0.8
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	9	0.8
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	11	0.8
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	12	0.8
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	14	0.8
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	19	0.8
(1,72)	1:159:A:GLN:H	1:160:A:ALA:H	8	0.8
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	2	0.8
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	3	0.8
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	14	0.8
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	6	0.8
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	3	0.8
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	7	0.8
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	10	0.8
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	4	0.8
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	20	0.79
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	17	0.79
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	20	0.79
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	17	0.79
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	7	0.79
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	1	0.79
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	10	0.79
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	5	0.79
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	17	0.79
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	12	0.79
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	19	0.79
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	4	0.79
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	5	0.79
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	10	0.79
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	15	0.79
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	16	0.79
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	18	0.79
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	19	0.79
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	20	0.79
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	4	0.79
(1,77)	1:97:A:GLN:H	1:99:A:CYS:H	4	0.79
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	10	0.79
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	11	0.79
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	16	0.79
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	11	0.79
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	20	0.79

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	17	0.79
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	14	0.79
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	3	0.79
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	1	0.79
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	3	0.79
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	4	0.78
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	7	0.78
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	1	0.78
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	1	0.78
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	6	0.78
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	6	0.78
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	6	0.78
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	6	0.78
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	6	0.78
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	6	0.78
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	6	0.78
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	6	0.78
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	6	0.78
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	4	0.78
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	7	0.78
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	1	0.78
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	1	0.78
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	19	0.78
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	11	0.78
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	8	0.78
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	11	0.78
(1,135)	1:64:A:CYS:H	1:66:A:PHE:H	7	0.78
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	3	0.78
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	14	0.78
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	6	0.78
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	14	0.78
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	10	0.78
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	3	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	1	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	3	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	5	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	8	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	9	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	13	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	14	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	17	0.78
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	19	0.78

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	15	0.78
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	19	0.78
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	7	0.78
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	10	0.78
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	16	0.78
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	18	0.78
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	14	0.78
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	12	0.78
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	19	0.78
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	7	0.77
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	14	0.77
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	14	0.77
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	14	0.77
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	14	0.77
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	14	0.77
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	14	0.77
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	14	0.77
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	14	0.77
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	14	0.77
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	7	0.77
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	19	0.77
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	7	0.77
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	15	0.77
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	13	0.77
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	15	0.77
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	10	0.77
(1,106)	1:96:A:SER:H	1:97:A:GLN:H	13	0.77
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	12	0.77
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	7	0.77
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	10	0.77
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	11	0.77
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	12	0.77
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	8	0.77
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	20	0.77
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	6	0.77
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	12	0.77
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	12	0.77
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	3	0.77
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	7	0.77
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	8	0.76
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	10	0.76
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	8	0.76

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	10	0.76
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	8	0.76
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	1	0.76
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	17	0.76
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	18	0.76
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	6	0.76
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	20	0.76
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	14	0.76
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	1	0.76
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	17	0.76
(1,47)	1:49:A:ASP:H	1:52:A:LEU:H	16	0.76
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	4	0.76
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	14	0.76
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	1	0.76
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	5	0.76
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	8	0.75
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	9	0.75
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	8	0.75
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	9	0.75
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	13	0.75
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	13	0.75
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	13	0.75
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	13	0.75
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	13	0.75
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	13	0.75
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	13	0.75
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	13	0.75
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	13	0.75
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	17	0.75
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	2	0.75
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	10	0.75
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	9	0.75
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	4	0.75
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	11	0.75
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	11	0.75
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	15	0.75
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	6	0.75
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	9	0.75
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	19	0.75
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	2	0.75
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	16	0.75
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	20	0.75

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,69)	1:42:A:ILE:H	1:126:A:ASP:H	15	0.75
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	3	0.75
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	14	0.75
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	4	0.75
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	20	0.75
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	2	0.75
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	9	0.75
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	14	0.75
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	16	0.75
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	3	0.75
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	10	0.75
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	16	0.75
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	20	0.74
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.74
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	18	0.74
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	20	0.74
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	20	0.74
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	18	0.74
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	10	0.74
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	13	0.74
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	8	0.74
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	18	0.74
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	3	0.74
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	7	0.74
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	18	0.74
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	1	0.74
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	10	0.74
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	14	0.74
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	2	0.74
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	15	0.74
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	4	0.74
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	7	0.74
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	2	0.74
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	10	0.74
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	7	0.74
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	11	0.74
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	6	0.74
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	6	0.74
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	8	0.74
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	11	0.74
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	3	0.74
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	8	0.74

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,146)	1:42:A:ILE:H	1:124:A:ASP:H	20	0.73
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	18	0.73
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	12	0.73
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	12	0.73
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	12	0.73
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	12	0.73
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	12	0.73
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	12	0.73
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	12	0.73
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	12	0.73
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	12	0.73
(3,28)	1:42:A:ILE:H	1:124:A:ASP:H	20	0.73
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	18	0.73
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	19	0.73
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	9	0.73
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	19	0.73
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	10	0.73
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	5	0.73
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	19	0.73
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	16	0.73
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	20	0.73
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	15	0.73
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	18	0.73
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	16	0.73
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	3	0.73
(1,206)	1:64:A:CYS:H	1:66:A:PHE:H	7	0.72
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	17	0.72
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	19	0.72
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	8	0.72
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	3	0.72
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	20	0.72
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	3	0.72
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	5	0.72
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	19	0.72
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	6	0.72
(1,96)	1:26:A:ILE:H	1:29:A:ASN:H	18	0.72
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	11	0.72
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	13	0.72
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	17	0.72
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	19	0.72
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	20	0.72
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	2	0.72

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	5	0.72
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	17	0.72
(1,24)	1:103:A:PHE:H	1:105:A:ARG:H	9	0.72
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	2	0.72
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	2	0.72
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	9	0.72
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	14	0.72
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	4	0.71
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	11	0.71
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	1	0.71
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	19	0.71
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	4	0.71
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	11	0.71
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	1	0.71
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	19	0.71
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	4	0.71
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	11	0.71
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	2	0.71
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	13	0.71
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	15	0.71
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	1	0.71
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	2	0.71
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	11	0.71
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	18	0.71
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	6	0.71
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	8	0.71
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	1	0.71
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	4	0.71
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	16	0.71
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	20	0.71
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	12	0.71
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	17	0.71
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	13	0.71
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	4	0.71
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	7	0.71
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	6	0.71
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	8	0.71
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	11	0.71
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	9	0.7
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	12	0.7
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	13	0.7
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	18	0.7

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	9	0.7
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	12	0.7
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	13	0.7
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	18	0.7
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	1	0.7
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	9	0.7
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	15	0.7
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	18	0.7
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	20	0.7
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	3	0.7
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	5	0.7
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	15	0.7
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	9	0.7
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	12	0.7
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	9	0.7
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	8	0.7
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	4	0.7
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	7	0.7
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	11	0.7
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	16	0.7
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	5	0.7
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	7	0.7
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	13	0.7
(1,79)	1:104:A:GLU:H	1:106:A:THR:H	18	0.7
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	15	0.7
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	4	0.7
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	10	0.7
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	1	0.7
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	16	0.69
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	16	0.69
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	16	0.69
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	16	0.69
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	2	0.69
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	7	0.69
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	11	0.69
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	11	0.69
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	1	0.69
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	17	0.69
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	5	0.69
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	6	0.69
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	13	0.69
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	14	0.69

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	18	0.69
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	9	0.69
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	18	0.69
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	17	0.69
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	18	0.69
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	5	0.69
(3,172)	1:137:A:SER:H	1:141:A:ILE:H	20	0.68
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	5	0.68
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	13	0.68
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	4	0.68
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	1	0.68
(3,54)	1:137:A:SER:H	1:141:A:ILE:H	20	0.68
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	5	0.68
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	13	0.68
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	4	0.68
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	1	0.68
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	1	0.68
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	9	0.68
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	8	0.68
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	20	0.68
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	9	0.68
(1,177)	1:54:A:HIS:H	1:56:A:SER:H	14	0.68
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	3	0.68
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	18	0.68
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	14	0.68
(1,137)	1:54:A:HIS:H	1:56:A:SER:H	14	0.68
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	13	0.68
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	15	0.68
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	11	0.68
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	16	0.68
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	18	0.68
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	15	0.68
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	4	0.68
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	13	0.68
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	15	0.68
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	6	0.68
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	13	0.68
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	1	0.67
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	13	0.67
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	12	0.67
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	13	0.67
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	18	0.67

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	12	0.67
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	12	0.67
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	12	0.67
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	12	0.67
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	12	0.67
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	12	0.67
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	12	0.67
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	12	0.67
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	12	0.67
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	1	0.67
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	13	0.67
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	12	0.67
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	13	0.67
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	18	0.67
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	9	0.67
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	15	0.67
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	19	0.67
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	14	0.67
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	13	0.67
(1,157)	1:40:A:HIS:H	1:41:A:VAL:H	18	0.67
(1,155)	1:30:A:SER:H	1:31:A:GLU:H	19	0.67
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	2	0.67
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	17	0.67
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	20	0.67
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	4	0.67
(1,70)	1:130:A:VAL:H	1:131:A:THR:H	6	0.67
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	5	0.67
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	9	0.67
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	19	0.67
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	6	0.67
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	18	0.67
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	12	0.67
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	11	0.67
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	12	0.66
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	3	0.66
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	15	0.66
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	19	0.66
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	12	0.66
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	3	0.66
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	15	0.66
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	19	0.66
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	11	0.66

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	20	0.66
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	10	0.66
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	3	0.66
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	17	0.66
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	16	0.66
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	8	0.66
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	17	0.66
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	17	0.66
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	12	0.66
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	19	0.66
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	3	0.66
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	15	0.66
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	4	0.66
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	8	0.66
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	18	0.66
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	11	0.66
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	18	0.66
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	19	0.66
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	3	0.65
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	14	0.65
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	3	0.65
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	7	0.65
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	4	0.65
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	16	0.65
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	16	0.65
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	16	0.65
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	16	0.65
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	16	0.65
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	16	0.65
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	16	0.65
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	16	0.65
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	16	0.65
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	3	0.65
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	14	0.65
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	3	0.65
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	7	0.65
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	4	0.65
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	4	0.65
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	18	0.65
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	14	0.65
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	13	0.65
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	4	0.65

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	3	0.65
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	13	0.65
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	4	0.65
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	5	0.65
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	6	0.65
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	9	0.65
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	16	0.65
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	6	0.65
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	10	0.65
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	17	0.65
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.65
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	13	0.65
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	14	0.65
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	4	0.65
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	7	0.64
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	11	0.64
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	20	0.64
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	17	0.64
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	7	0.64
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	11	0.64
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	20	0.64
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	17	0.64
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	3	0.64
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	15	0.64
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	4	0.64
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	13	0.64
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	17	0.64
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	15	0.64
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	10	0.64
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	15	0.64
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	9	0.64
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	20	0.64
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	9	0.64
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	1	0.64
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	11	0.64
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	7	0.63
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	10	0.63
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	16	0.63
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	14	0.63
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	12	0.63
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	19	0.63
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	7	0.63

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	10	0.63
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	16	0.63
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	14	0.63
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	12	0.63
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	19	0.63
(1,209)	1:21:A:TRP:H	1:87:A:SER:H	4	0.63
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	4	0.63
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	13	0.63
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	2	0.63
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	17	0.63
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	5	0.63
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	3	0.63
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	11	0.63
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	13	0.63
(1,150)	1:58:A:THR:H	1:60:A:LEU:H	3	0.63
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	15	0.63
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	6	0.63
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	3	0.63
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	17	0.63
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	4	0.63
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	19	0.63
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	12	0.63
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	5	0.63
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	1	0.63
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	7	0.62
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	10	0.62
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	6	0.62
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	3	0.62
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	7	0.62
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	7	0.62
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	10	0.62
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	6	0.62
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	3	0.62
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	7	0.62
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	17	0.62
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	2	0.62
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	8	0.62
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	16	0.62
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	18	0.62
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	9	0.62
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	10	0.62
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	14	0.62

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	15	0.62
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	5	0.62
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	19	0.62
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	2	0.62
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	1	0.62
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	2	0.62
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	16	0.62
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	18	0.62
(1,160)	1:46:A:THR:H	1:47:A:LYS:H	20	0.62
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	12	0.62
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	14	0.62
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	15	0.62
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	9	0.62
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	19	0.62
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	18	0.62
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	17	0.62
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.62
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	1	0.61
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	17	0.61
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	4	0.61
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	14	0.61
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	5	0.61
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	1	0.61
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	17	0.61
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	4	0.61
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	14	0.61
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	5	0.61
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	15	0.61
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	10	0.61
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	13	0.61
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	2	0.61
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	4	0.61
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	18	0.61
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	15	0.61
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	16	0.61
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	6	0.61
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	4	0.61
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	14	0.61
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	15	0.61
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	11	0.61
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	15	0.61
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	17	0.61

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	20	0.61
(1,53)	1:30:A:SER:H	1:31:A:GLU:H	19	0.61
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	10	0.61
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	2	0.6
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	10	0.6
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	6	0.6
(3,142)	1:42:A:ILE:H	1:126:A:ASP:H	15	0.6
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	3	0.6
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	14	0.6
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	2	0.6
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	9	0.6
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	14	0.6
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	3	0.6
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	10	0.6
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	16	0.6
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	13	0.6
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	13	0.6
(3,60)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	13	0.6
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	13	0.6
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	13	0.6
(3,60)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	13	0.6
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	13	0.6
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	13	0.6
(3,60)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	13	0.6
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	2	0.6
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	10	0.6
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	6	0.6
(3,24)	1:42:A:ILE:H	1:126:A:ASP:H	15	0.6
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	3	0.6
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	14	0.6
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	2	0.6
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	9	0.6
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	14	0.6
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	3	0.6
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	10	0.6
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	16	0.6
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	17	0.6
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	5	0.6
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	16	0.6
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	17	0.6
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	19	0.6
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	20	0.6

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	1	0.6
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	8	0.6
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	6	0.6
(1,129)	1:180:A:TYR:H	1:181:A:LEU:H	6	0.6
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	3	0.6
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	15	0.6
(1,105)	1:67:A:ALA:H	1:68:A:LYS:H	19	0.6
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	4	0.6
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	19	0.6
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	1	0.6
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	6	0.6
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	2	0.6
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	10	0.6
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	12	0.6
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	13	0.6
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	20	0.6
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	17	0.6
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	5	0.6
(1,15)	1:180:A:TYR:H	1:181:A:LEU:H	6	0.6
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	6	0.6
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	5	0.6
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	10	0.59
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	18	0.59
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	1	0.59
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	2	0.59
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	2	0.59
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	10	0.59
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	6	0.59
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	8	0.59
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	11	0.59
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	10	0.59
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	18	0.59
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	1	0.59
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	2	0.59
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	2	0.59
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	10	0.59
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	6	0.59
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	8	0.59
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	11	0.59
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	16	0.59
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	16	0.59
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	16	0.59

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	16	0.59
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	16	0.59
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	16	0.59
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	16	0.59
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	16	0.59
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	16	0.59
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	2	0.59
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	5	0.59
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	6	0.59
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	7	0.59
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	14	0.59
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	3	0.59
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	10	0.59
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	12	0.59
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	12	0.59
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	5	0.59
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	11	0.59
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	13	0.59
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	2	0.59
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	8	0.59
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	16	0.59
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	18	0.59
(1,68)	1:41:A:VAL:H	1:126:A:ASP:H	11	0.59
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	7	0.59
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	13	0.59
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	5	0.59
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	10	0.59
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	12	0.59
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	14	0.59
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	15	0.58
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	3	0.58
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	15	0.58
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	3	0.58
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	12	0.58
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	16	0.58
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	12	0.58
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	3	0.58
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	11	0.58
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	1	0.58
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	4	0.58
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	3	0.58
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	7	0.58

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,149)	1:68:A:LYS:H	1:85:A:ARG:H	4	0.58
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	16	0.58
(1,118)	1:68:A:LYS:H	1:85:A:ARG:H	4	0.58
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	13	0.58
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	10	0.58
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	13	0.58
(1,74)	1:60:A:LEU:H	1:62:A:ASN:H	8	0.58
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	12	0.58
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	1	0.58
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	16	0.58
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	3	0.58
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	5	0.58
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	8	0.58
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	13	0.58
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	16	0.58
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	17	0.58
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	18	0.58
(1,36)	1:73:A:TYR:H	1:81:A:TYR:H	15	0.58
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	19	0.57
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	20	0.57
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	2	0.57
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	5	0.57
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	2	0.57
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	9	0.57
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	14	0.57
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	19	0.57
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	20	0.57
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	2	0.57
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	5	0.57
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	2	0.57
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	9	0.57
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	14	0.57
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	7	0.57
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	15	0.57
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	6	0.57
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	14	0.57
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	15	0.57
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	1	0.57
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	2	0.57
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	19	0.57
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	18	0.57
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	2	0.57

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	6	0.57
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	7	0.57
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	11	0.57
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	15	0.57
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	19	0.57
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	20	0.57
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	7	0.57
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	2	0.56
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	11	0.56
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	16	0.56
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	13	0.56
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	6	0.56
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	8	0.56
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	11	0.56
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	2	0.56
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	11	0.56
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	16	0.56
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	13	0.56
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	6	0.56
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	8	0.56
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	11	0.56
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	5	0.56
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	7	0.56
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	1	0.56
(1,197)	1:123:A:GLY:H	1:124:A:ASP:H	12	0.56
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	20	0.56
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	7	0.56
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	8	0.56
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	13	0.56
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	6	0.56
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	9	0.56
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	5	0.55
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	9	0.55
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	15	0.55
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	10	0.55
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	5	0.55
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	9	0.55
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	15	0.55
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	10	0.55
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	8	0.55
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	4	0.55
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	15	0.55

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	6	0.55
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	20	0.55
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	18	0.55
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	11	0.55
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	10	0.55
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	10	0.55
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	2	0.55
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	10	0.55
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	17	0.55
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	12	0.55
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	3	0.55
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	8	0.55
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	1	0.55
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	9	0.55
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	2	0.55
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	16	0.55
(1,5)	1:73:A:TYR:H	1:81:A:TYR:H	15	0.55
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	18	0.54
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	5	0.54
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	18	0.54
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	5	0.54
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	6	0.54
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	11	0.54
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	16	0.54
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	20	0.54
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	19	0.54
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	20	0.54
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	15	0.54
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	6	0.54
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	6	0.54
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	12	0.54
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	14	0.54
(1,37)	1:72:A:PHE:H	1:73:A:TYR:H	4	0.54
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	19	0.54
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	1	0.53
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	9	0.53
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	9	0.53
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	6	0.53
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	13	0.53
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	1	0.53
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	9	0.53
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	9	0.53

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	6	0.53
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	13	0.53
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	7	0.53
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	6	0.53
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	16	0.53
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	20	0.53
(1,131)	1:99:A:CYS:H	1:101:A:SER:H	18	0.53
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	18	0.53
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	5	0.53
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	7	0.53
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	1	0.53
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	7	0.53
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	9	0.52
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	19	0.52
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	11	0.52
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	9	0.52
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	19	0.52
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	11	0.52
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	5	0.52
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	19	0.52
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	12	0.52
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	8	0.52
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	15	0.52
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	15	0.52
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	7	0.52
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	12	0.52
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	11	0.51
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	20	0.51
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	17	0.51
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	3	0.51
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	15	0.51
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	18	0.51
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	18	0.51
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	19	0.51
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	11	0.51
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	20	0.51
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	17	0.51
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	3	0.51
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	15	0.51
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	18	0.51
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	18	0.51
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	19	0.51

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	17	0.51
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	13	0.51
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	14	0.51
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	12	0.51
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	11	0.51
(1,161)	1:46:A:THR:H	1:48:A:GLN:H	9	0.51
(1,115)	1:179:A:LYS:H	1:181:A:LEU:H	19	0.51
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	6	0.51
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	14	0.51
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	19	0.51
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	9	0.51
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	4	0.5
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	18	0.5
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	13	0.5
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	4	0.5
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	10	0.5
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	17	0.5
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.5
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	4	0.5
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	4	0.5
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	18	0.5
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	13	0.5
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	4	0.5
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	10	0.5
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	17	0.5
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.5
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	4	0.5
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	10	0.5
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	6	0.5
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	7	0.5
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	12	0.5
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	3	0.5
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	4	0.5
(1,19)	1:76:A:CYS:H	1:78:A:TYR:H	17	0.5
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	17	0.5
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	3	0.49
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	17	0.49
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	15	0.49
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	3	0.49
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	17	0.49
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	15	0.49
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	12	0.49

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	12	0.49
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	12	0.49
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	12	0.49
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	12	0.49
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	12	0.49
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	12	0.49
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	12	0.49
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	12	0.49
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	3	0.49
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	3	0.49
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	3	0.49
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	3	0.49
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	3	0.49
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	3	0.49
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	3	0.49
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	3	0.49
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	3	0.49
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	14	0.49
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	13	0.49
(1,167)	1:104:A:GLU:H	1:106:A:THR:H	18	0.49
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	4	0.49
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	4	0.49
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	17	0.49
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	5	0.49
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	6	0.49
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	12	0.49
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	18	0.49
(1,58)	1:93:A:VAL:H	1:155:A:LEU:H	2	0.49
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	3	0.49
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	18	0.49
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	9	0.49
(1,3)	1:119:A:LEU:H	1:163:A:PHE:H	19	0.49
(3,170)	1:21:A:TRP:H	1:87:A:SER:H	4	0.48
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	3	0.48
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	11	0.48
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	13	0.48
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	19	0.48
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	12	0.48
(3,52)	1:21:A:TRP:H	1:87:A:SER:H	4	0.48
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	3	0.48
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	11	0.48
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	13	0.48

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	19	0.48
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	12	0.48
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	9	0.48
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	14	0.48
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	9	0.48
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	19	0.48
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	12	0.48
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	10	0.48
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	2	0.48
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	9	0.48
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	5	0.48
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	13	0.48
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	2	0.48
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	8	0.48
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	16	0.48
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	1	0.48
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	13	0.48
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	17	0.48
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	1	0.47
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	2	0.47
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	16	0.47
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	18	0.47
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	17	0.47
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.47
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	1	0.47
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	2	0.47
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	16	0.47
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	18	0.47
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	17	0.47
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	20	0.47
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	12	0.47
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	14	0.47
(1,162)	1:46:A:THR:H	1:50:A:LEU:H	20	0.47
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	4	0.47
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	14	0.47
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	15	0.47
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	20	0.47
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	20	0.47
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	15	0.46
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	15	0.46
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	4	0.46
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	14	0.46

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	15	0.46
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	15	0.46
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	4	0.46
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	14	0.46
(1,204)	1:48:A:GLN:H	1:50:A:LEU:H	1	0.46
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	8	0.46
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	17	0.46
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	11	0.46
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	15	0.46
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	17	0.46
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	9	0.46
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	2	0.46
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	9	0.46
(1,155)	1:30:A:SER:H	1:31:A:GLU:H	14	0.46
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	9	0.46
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	12	0.46
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	2	0.46
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	12	0.46
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	1	0.46
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	20	0.46
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	19	0.46
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	9	0.46
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	10	0.46
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	12	0.46
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	17	0.45
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	17	0.45
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	6	0.45
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	17	0.45
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	17	0.45
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	6	0.45
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	19	0.45
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	5	0.45
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	1	0.45
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	13	0.45
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	9	0.45
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	5	0.45
(1,60)	1:27:A:GLU:H	1:29:A:ASN:H	3	0.45
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	5	0.45
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	5	0.45
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	19	0.45
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	10	0.45
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	8	0.45

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	2	0.44
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	5	0.44
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	12	0.44
(3,141)	1:41:A:VAL:H	1:126:A:ASP:H	11	0.44
(3,65)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	16	0.44
(3,65)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	16	0.44
(3,65)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	16	0.44
(3,65)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	16	0.44
(3,65)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	16	0.44
(3,65)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	16	0.44
(3,65)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	16	0.44
(3,65)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	16	0.44
(3,65)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	16	0.44
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	2	0.44
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	5	0.44
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	12	0.44
(3,23)	1:41:A:VAL:H	1:126:A:ASP:H	11	0.44
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	1	0.44
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	7	0.44
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	13	0.44
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	8	0.44
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	5	0.44
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	14	0.44
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	17	0.44
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	14	0.44
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	16	0.44
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	16	0.44
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	7	0.44
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	12	0.43
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	16	0.43
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	7	0.43
(3,158)	1:68:A:LYS:H	1:85:A:ARG:H	4	0.43
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	16	0.43
(3,154)	1:68:A:LYS:H	1:85:A:ARG:H	4	0.43
(3,135)	1:73:A:TYR:H	1:81:A:TYR:H	15	0.43
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	12	0.43
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	16	0.43
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	7	0.43
(3,40)	1:68:A:LYS:H	1:85:A:ARG:H	4	0.43
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	16	0.43
(3,36)	1:68:A:LYS:H	1:85:A:ARG:H	4	0.43
(3,17)	1:73:A:TYR:H	1:81:A:TYR:H	15	0.43

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	3	0.43
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	3	0.43
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	3	0.43
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	3	0.43
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	3	0.43
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	3	0.43
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	3	0.43
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	3	0.43
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	3	0.43
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	8	0.43
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	19	0.43
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	12	0.43
(1,165)	1:137:A:SER:H	1:139:A:PHE:H	14	0.43
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	7	0.43
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	5	0.43
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	6	0.43
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	14	0.43
(1,98)	1:48:A:GLN:H	1:50:A:LEU:H	1	0.43
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	1	0.43
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	7	0.43
(1,57)	1:87:A:SER:H	1:89:A:HIS:H	8	0.43
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	17	0.43
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	2	0.43
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	19	0.43
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	16	0.43
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	6	0.42
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	14	0.42
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	15	0.42
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	2	0.42
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	7	0.42
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	6	0.42
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	14	0.42
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	15	0.42
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	2	0.42
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	7	0.42
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	4	0.42
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	15	0.42
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	17	0.42
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	18	0.42
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	18	0.42
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	1	0.42
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	10	0.42

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	14	0.42
(1,64)	1:86:A:ASP:H	1:87:A:SER:H	19	0.42
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	13	0.42
(1,16)	1:179:A:LYS:H	1:181:A:LEU:H	19	0.42
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	8	0.42
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	8	0.41
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	13	0.41
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	8	0.41
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	13	0.41
(1,266)	1:18:A:LEU:HD21	1:119:A:LEU:HD11	8	0.41
(1,266)	1:18:A:LEU:HD21	1:119:A:LEU:HD12	8	0.41
(1,266)	1:18:A:LEU:HD21	1:119:A:LEU:HD13	8	0.41
(1,266)	1:18:A:LEU:HD22	1:119:A:LEU:HD11	8	0.41
(1,266)	1:18:A:LEU:HD22	1:119:A:LEU:HD12	8	0.41
(1,266)	1:18:A:LEU:HD22	1:119:A:LEU:HD13	8	0.41
(1,266)	1:18:A:LEU:HD23	1:119:A:LEU:HD11	8	0.41
(1,266)	1:18:A:LEU:HD23	1:119:A:LEU:HD12	8	0.41
(1,266)	1:18:A:LEU:HD23	1:119:A:LEU:HD13	8	0.41
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	1	0.41
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	14	0.41
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	20	0.41
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	4	0.41
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	11	0.41
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	4	0.41
(1,107)	1:67:A:ALA:H	1:120:A:MET:H	12	0.41
(1,39)	1:41:A:VAL:H	1:125:A:ALA:H	6	0.41
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	8	0.41
(1,3)	1:119:A:LEU:H	1:163:A:PHE:H	14	0.41
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	8	0.4
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	10	0.4
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	10	0.4
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	2	0.4
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	8	0.4
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	16	0.4
(3,122)	1:73:A:TYR:H	1:81:A:TYR:H	15	0.4
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	8	0.4
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	10	0.4
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	10	0.4
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	2	0.4
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	8	0.4
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	16	0.4
(3,4)	1:73:A:TYR:H	1:81:A:TYR:H	15	0.4

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	11	0.4
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	19	0.4
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	3	0.4
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	13	0.4
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	17	0.4
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	16	0.4
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	8	0.4
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	12	0.4
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	9	0.4
(1,53)	1:30:A:SER:H	1:31:A:GLU:H	14	0.4
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	9	0.4
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	1	0.4
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	3	0.4
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	3	0.4
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	4	0.4
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	12	0.4
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	15	0.39
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	19	0.39
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	15	0.39
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	19	0.39
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD11	13	0.39
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD12	13	0.39
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD13	13	0.39
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD11	13	0.39
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD12	13	0.39
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD13	13	0.39
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD11	13	0.39
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD12	13	0.39
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD13	13	0.39
(1,209)	1:21:A:TRP:H	1:87:A:SER:H	8	0.39
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	5	0.39
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	7	0.39
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	6	0.39
(1,166)	1:137:A:SER:H	1:140:A:LYS:H	20	0.39
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	5	0.39
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	14	0.39
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	7	0.39
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	5	0.39
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	20	0.39
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	3	0.39
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	10	0.39
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	17	0.39

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	9	0.39
(1,11)	1:143:A:LEU:H	1:145:A:ASN:H	15	0.39
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	1	0.38
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	1	0.38
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	15	0.38
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	20	0.38
(1,209)	1:21:A:TRP:H	1:87:A:SER:H	20	0.38
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	13	0.38
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	17	0.38
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	19	0.38
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	10	0.38
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	18	0.38
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	11	0.38
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	14	0.38
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	4	0.38
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	5	0.37
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	19	0.37
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	8	0.37
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	15	0.37
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	12	0.37
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	5	0.37
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	19	0.37
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	8	0.37
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	15	0.37
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	12	0.37
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD11	11	0.37
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD12	11	0.37
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD13	11	0.37
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD11	11	0.37
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD12	11	0.37
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD13	11	0.37
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD11	11	0.37
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD12	11	0.37
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD13	11	0.37
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	20	0.37
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	12	0.37
(1,178)	1:90:A:GLN:H	1:177:A:VAL:H	7	0.37
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	2	0.37
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	10	0.37
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	16	0.37
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	17	0.37
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	16	0.37

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	20	0.37
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	12	0.37
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	4	0.37
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	11	0.37
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	9	0.37
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	17	0.36
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	13	0.36
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	14	0.36
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	17	0.36
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	13	0.36
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	14	0.36
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	6	0.36
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	14	0.36
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	10	0.36
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	9	0.36
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	5	0.36
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	3	0.36
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	8	0.36
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	9	0.36
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	15	0.36
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	17	0.36
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	7	0.36
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	1	0.36
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	11	0.36
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	5	0.36
(1,3)	1:119:A:LEU:H	1:163:A:PHE:H	3	0.36
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	4	0.36
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	10	0.35
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	17	0.35
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	10	0.35
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	17	0.35
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG11	19	0.35
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG12	19	0.35
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG13	19	0.35
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG11	19	0.35
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG12	19	0.35
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG13	19	0.35
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG11	19	0.35
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG12	19	0.35
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG13	19	0.35
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	13	0.35
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	13	0.35

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	13	0.35
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	13	0.35
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	13	0.35
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	13	0.35
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	13	0.35
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	13	0.35
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	13	0.35
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	11	0.35
(1,178)	1:90:A:GLN:H	1:177:A:VAL:H	4	0.35
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	15	0.35
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	18	0.35
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	9	0.35
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	5	0.35
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	19	0.35
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	1	0.35
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	2	0.35
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	3	0.35
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	5	0.35
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	15	0.35
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	17	0.35
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	19	0.35
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	20	0.35
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	14	0.34
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	4	0.34
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	17	0.34
(3,139)	1:93:A:VAL:H	1:155:A:LEU:H	2	0.34
(3,120)	1:119:A:LEU:H	1:163:A:PHE:H	19	0.34
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	12	0.34
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	12	0.34
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	12	0.34
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	12	0.34
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	12	0.34
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	12	0.34
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	12	0.34
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	12	0.34
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	12	0.34
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	3	0.34
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	3	0.34
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	3	0.34
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	3	0.34
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	3	0.34
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	3	0.34

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	3	0.34
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	3	0.34
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	3	0.34
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	14	0.34
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	4	0.34
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	17	0.34
(3,21)	1:93:A:VAL:H	1:155:A:LEU:H	2	0.34
(3,2)	1:119:A:LEU:H	1:163:A:PHE:H	19	0.34
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	1	0.34
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	1	0.34
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	1	0.34
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	1	0.34
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	1	0.34
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	1	0.34
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	1	0.34
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	1	0.34
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	1	0.34
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	6	0.34
(1,213)	1:136:A:MET:H	1:141:A:ILE:H	19	0.34
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	3	0.34
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	16	0.34
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	5	0.34
(1,185)	1:138:A:HIS:H	1:140:A:LYS:H	18	0.34
(1,178)	1:90:A:GLN:H	1:177:A:VAL:H	20	0.34
(1,163)	1:135:A:ALA:H	1:137:A:SER:H	8	0.34
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	1	0.34
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	9	0.34
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	19	0.34
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	2	0.34
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	9	0.34
(1,110)	1:45:A:LEU:H	1:124:A:ASP:H	7	0.34
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	2	0.34
(1,83)	1:138:A:HIS:H	1:140:A:LYS:H	18	0.34
(1,17)	1:119:A:LEU:H	1:163:A:PHE:H	16	0.34
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	18	0.34
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	7	0.34
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	8	0.34
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	11	0.34
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	14	0.34
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	18	0.34
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	9	0.33
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	19	0.33

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	12	0.33
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	10	0.33
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	5	0.33
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	9	0.33
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	19	0.33
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	12	0.33
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	10	0.33
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	5	0.33
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	9	0.33
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	2	0.33
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	8	0.33
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	16	0.33
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	15	0.33
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	16	0.33
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	8	0.33
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	17	0.33
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	14	0.33
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	3	0.33
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	11	0.33
(1,88)	1:42:A:ILE:H	1:125:A:ALA:H	1	0.33
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	3	0.33
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	8	0.33
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	13	0.33
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	16	0.33
(1,41)	1:42:A:ILE:H	1:125:A:ALA:H	1	0.33
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	19	0.33
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	6	0.33
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	16	0.33
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	14	0.33
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	6	0.33
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	10	0.33
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	13	0.33
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	16	0.33
(3,162)	1:46:A:THR:H	1:50:A:LEU:H	20	0.32
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	4	0.32
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	15	0.32
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	20	0.32
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	20	0.32
(3,44)	1:46:A:THR:H	1:50:A:LEU:H	20	0.32
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	4	0.32
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	15	0.32
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	20	0.32

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	20	0.32
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG21	10	0.32
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG22	10	0.32
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG23	10	0.32
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG21	10	0.32
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG22	10	0.32
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG23	10	0.32
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG21	10	0.32
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG22	10	0.32
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG23	10	0.32
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	17	0.32
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	4	0.32
(1,155)	1:30:A:SER:H	1:31:A:GLU:H	13	0.32
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	5	0.32
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	6	0.32
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	17	0.32
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	18	0.32
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	7	0.32
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	1	0.32
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	4	0.32
(1,126)	1:86:A:ASP:H	1:88:A:SER:H	4	0.32
(1,126)	1:86:A:ASP:H	1:88:A:SER:H	20	0.32
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	3	0.32
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	2	0.32
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	7	0.32
(1,3)	1:119:A:LEU:H	1:163:A:PHE:H	18	0.32
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	17	0.31
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	9	0.31
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	12	0.31
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	19	0.31
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	17	0.31
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	9	0.31
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	12	0.31
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	19	0.31
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG21	16	0.31
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG22	16	0.31
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG23	16	0.31
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG21	16	0.31
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG22	16	0.31
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG23	16	0.31
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG21	16	0.31
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG22	16	0.31

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG23	16	0.31
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	9	0.31
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	9	0.31
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	9	0.31
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	10	0.31
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	2	0.31
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	20	0.31
(1,144)	1:44:A:SER:H	1:124:A:ASP:H	18	0.31
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	17	0.31
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	6	0.31
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	5	0.31
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	8	0.31
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	13	0.31
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	19	0.31
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	13	0.31
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	2	0.31
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	10	0.31
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	12	0.31
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	5	0.31
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	13	0.31
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	12	0.31
(1,2)	1:82:A:ALA:H	1:83:A:SER:H	12	0.31
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	19	0.3
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	5	0.3
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	13	0.3
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	5	0.3
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	10	0.3
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	19	0.3
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	5	0.3
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	13	0.3
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	5	0.3
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	10	0.3
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	14	0.3
(1,180)	1:90:A:GLN:H	1:178:A:TYR:H	16	0.3
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	5	0.3
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	2	0.3
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	7	0.3
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	15	0.3
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	19	0.3
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	13	0.3
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	5	0.3
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	18	0.3

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	4	0.3
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	16	0.3
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	3	0.3
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	18	0.3
(1,3)	1:119:A:LEU:H	1:163:A:PHE:H	11	0.3
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	1	0.29
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	14	0.29
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	17	0.29
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	1	0.29
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	14	0.29
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	17	0.29
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	13	0.29
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	13	0.29
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	13	0.29
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	13	0.29
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	13	0.29
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	13	0.29
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	13	0.29
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	13	0.29
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	13	0.29
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	14	0.29
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	3	0.29
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	11	0.29
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	13	0.29
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	5	0.29
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	3	0.29
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	1	0.29
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	3	0.29
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	10	0.29
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	12	0.29
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	4	0.29
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	4	0.29
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	12	0.29
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	11	0.29
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	15	0.29
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	20	0.29
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	3	0.29
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	11	0.29
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	13	0.29
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	8	0.28
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	19	0.28
(3,134)	1:131:A:THR:H	1:135:A:ALA:H	17	0.28

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	2	0.28
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	19	0.28
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	3	0.28
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	3	0.28
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	3	0.28
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	3	0.28
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	3	0.28
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	3	0.28
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	3	0.28
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	3	0.28
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	3	0.28
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	8	0.28
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	19	0.28
(3,16)	1:131:A:THR:H	1:135:A:ALA:H	17	0.28
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	2	0.28
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	19	0.28
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD11	4	0.28
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD12	4	0.28
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD13	4	0.28
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD11	4	0.28
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD12	4	0.28
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD13	4	0.28
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD11	4	0.28
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD12	4	0.28
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD13	4	0.28
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	9	0.28
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	9	0.28
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	9	0.28
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	9	0.28
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	9	0.28
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	9	0.28
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	9	0.28
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	9	0.28
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	9	0.28
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	1	0.28
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	2	0.28
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	16	0.28
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	18	0.28
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	2	0.28
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	20	0.28
(1,126)	1:86:A:ASP:H	1:88:A:SER:H	12	0.28
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	7	0.28

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	4	0.28
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	11	0.28
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	1	0.28
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	10	0.28
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	14	0.28
(1,42)	1:128:A:VAL:H	1:154:A:VAL:H	18	0.28
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	14	0.28
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	13	0.27
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	8	0.27
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	13	0.27
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	8	0.27
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	1	0.27
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	1	0.27
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	1	0.27
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	1	0.27
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	1	0.27
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	1	0.27
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	1	0.27
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	1	0.27
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	1	0.27
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD11	9	0.27
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD12	9	0.27
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD13	9	0.27
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD11	9	0.27
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD12	9	0.27
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD13	9	0.27
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD11	9	0.27
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD12	9	0.27
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD13	9	0.27
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	11	0.27
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	17	0.27
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	5	0.27
(1,192)	1:86:A:ASP:H	1:87:A:SER:H	19	0.27
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	16	0.27
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	4	0.27
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	11	0.27
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	13	0.27
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	3	0.27
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	13	0.27
(1,107)	1:67:A:ALA:H	1:120:A:MET:H	1	0.27
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	9	0.27
(1,94)	1:121:A:ILE:H	1:163:A:PHE:H	5	0.27

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	8	0.27
(1,61)	1:71:A:LYS:H	1:72:A:PHE:H	6	0.27
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	8	0.27
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	11	0.26
(3,150)	1:67:A:ALA:H	1:120:A:MET:H	12	0.26
(3,136)	1:41:A:VAL:H	1:125:A:ALA:H	6	0.26
(3,120)	1:119:A:LEU:H	1:163:A:PHE:H	14	0.26
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	11	0.26
(3,32)	1:67:A:ALA:H	1:120:A:MET:H	12	0.26
(3,18)	1:41:A:VAL:H	1:125:A:ALA:H	6	0.26
(3,2)	1:119:A:LEU:H	1:163:A:PHE:H	14	0.26
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	7	0.26
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	15	0.26
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	13	0.26
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	19	0.26
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	10	0.26
(1,152)	1:41:A:VAL:H	1:125:A:ALA:H	6	0.26
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	14	0.26
(1,126)	1:86:A:ASP:H	1:88:A:SER:H	17	0.26
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	20	0.26
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	6	0.26
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	11	0.26
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	12	0.26
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	2	0.26
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	18	0.26
(1,53)	1:30:A:SER:H	1:31:A:GLU:H	13	0.26
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	4	0.26
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	2	0.26
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	14	0.26
(1,7)	1:75:A:VAL:H	1:81:A:TYR:H	14	0.26
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	8	0.25
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	12	0.25
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	9	0.25
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	1	0.25
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	3	0.25
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	8	0.25
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	12	0.25
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	9	0.25
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	1	0.25
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	3	0.25
(1,271)	1:18:A:LEU:HD11	1:119:A:LEU:HD11	11	0.25
(1,271)	1:18:A:LEU:HD11	1:119:A:LEU:HD12	11	0.25

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,271)	1:18:A:LEU:HD11	1:119:A:LEU:HD13	11	0.25
(1,271)	1:18:A:LEU:HD12	1:119:A:LEU:HD11	11	0.25
(1,271)	1:18:A:LEU:HD12	1:119:A:LEU:HD12	11	0.25
(1,271)	1:18:A:LEU:HD12	1:119:A:LEU:HD13	11	0.25
(1,271)	1:18:A:LEU:HD13	1:119:A:LEU:HD11	11	0.25
(1,271)	1:18:A:LEU:HD13	1:119:A:LEU:HD12	11	0.25
(1,271)	1:18:A:LEU:HD13	1:119:A:LEU:HD13	11	0.25
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	12	0.25
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	4	0.25
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	11	0.25
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	3	0.25
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	10	0.25
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	12	0.25
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	5	0.25
(1,122)	1:151:A:ILE:H	1:152:A:VAL:H	7	0.25
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	8	0.25
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	9	0.25
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	5	0.25
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	15	0.25
(3,170)	1:21:A:TRP:H	1:87:A:SER:H	8	0.24
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	5	0.24
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD11	13	0.24
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD12	13	0.24
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD13	13	0.24
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD11	13	0.24
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD12	13	0.24
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD13	13	0.24
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD11	13	0.24
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD12	13	0.24
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD13	13	0.24
(3,52)	1:21:A:TRP:H	1:87:A:SER:H	8	0.24
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	5	0.24
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	4	0.24
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	4	0.24
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	4	0.24
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	4	0.24
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	4	0.24
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	4	0.24
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	4	0.24
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	4	0.24
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	4	0.24
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	10	0.24

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	10	0.24
(1,253)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	10	0.24
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	10	0.24
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	10	0.24
(1,253)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	10	0.24
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	10	0.24
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	10	0.24
(1,253)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	10	0.24
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	12	0.24
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	12	0.24
(1,224)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	12	0.24
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	12	0.24
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	12	0.24
(1,224)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	12	0.24
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	12	0.24
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	12	0.24
(1,224)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	12	0.24
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	6	0.24
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	16	0.24
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	7	0.24
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	13	0.24
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	4	0.24
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	19	0.24
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	9	0.24
(1,125)	1:88:A:SER:H	1:89:A:HIS:H	15	0.24
(1,110)	1:45:A:LEU:H	1:124:A:ASP:H	5	0.24
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	4	0.24
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	14	0.24
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	17	0.24
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	6	0.24
(1,26)	1:75:A:VAL:H	1:76:A:CYS:H	8	0.24
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	15	0.23
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	20	0.23
(3,170)	1:21:A:TRP:H	1:87:A:SER:H	20	0.23
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	14	0.23
(3,126)	1:126:A:ASP:H	1:156:A:ASN:H	4	0.23
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	15	0.23
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	20	0.23
(3,52)	1:21:A:TRP:H	1:87:A:SER:H	20	0.23
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	14	0.23
(3,8)	1:126:A:ASP:H	1:156:A:ASN:H	4	0.23
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	12	0.23

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	6	0.23
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	14	0.23
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	15	0.23
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	8	0.23
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	14	0.23
(1,126)	1:86:A:ASP:H	1:88:A:SER:H	11	0.23
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	18	0.23
(1,107)	1:67:A:ALA:H	1:120:A:MET:H	6	0.23
(1,94)	1:121:A:ILE:H	1:163:A:PHE:H	1	0.23
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	5	0.23
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	3	0.23
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	11	0.23
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	11	0.23
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	16	0.23
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	12	0.22
(3,164)	1:90:A:GLN:H	1:177:A:VAL:H	7	0.22
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	12	0.22
(3,46)	1:90:A:GLN:H	1:177:A:VAL:H	7	0.22
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG21	2	0.22
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG22	2	0.22
(1,273)	1:130:A:VAL:HG11	1:152:A:VAL:HG23	2	0.22
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG21	2	0.22
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG22	2	0.22
(1,273)	1:130:A:VAL:HG12	1:152:A:VAL:HG23	2	0.22
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG21	2	0.22
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG22	2	0.22
(1,273)	1:130:A:VAL:HG13	1:152:A:VAL:HG23	2	0.22
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	3	0.22
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	4	0.22
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	10	0.22
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	3	0.22
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	8	0.22
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	10	0.22
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	19	0.22
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	13	0.22
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	18	0.22
(1,125)	1:88:A:SER:H	1:89:A:HIS:H	11	0.22
(1,113)	1:124:A:ASP:H	1:159:A:GLN:H	8	0.22
(1,104)	1:84:A:ILE:H	1:91:A:ILE:H	20	0.22
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	10	0.22
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	4	0.22
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	10	0.22

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,21)	1:128:A:VAL:H	1:154:A:VAL:H	18	0.22
(3,143)	1:130:A:VAL:H	1:152:A:VAL:H	7	0.21
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	11	0.21
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	5	0.21
(3,120)	1:119:A:LEU:H	1:163:A:PHE:H	3	0.21
(3,25)	1:130:A:VAL:H	1:152:A:VAL:H	7	0.21
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	11	0.21
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	5	0.21
(3,2)	1:119:A:LEU:H	1:163:A:PHE:H	3	0.21
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	1	0.21
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	1	0.21
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	1	0.21
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	1	0.21
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	1	0.21
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	1	0.21
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	1	0.21
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	1	0.21
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	1	0.21
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	3	0.21
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	18	0.21
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	2	0.21
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	9	0.21
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	11	0.21
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	16	0.21
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	10	0.21
(1,201)	1:47:A:LYS:H	1:49:A:ASP:H	2	0.21
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	11	0.21
(1,185)	1:138:A:HIS:H	1:140:A:LYS:H	2	0.21
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	18	0.21
(1,159)	1:46:A:THR:H	1:49:A:ASP:H	10	0.21
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	8	0.21
(1,84)	1:35:A:TYR:H	1:36:A:HIS:H	16	0.21
(1,83)	1:138:A:HIS:H	1:140:A:LYS:H	2	0.21
(1,55)	1:88:A:SER:H	1:89:A:HIS:H	15	0.21
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	1	0.21
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	9	0.21
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	13	0.21
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	10	0.21
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	6	0.21
(1,18)	1:119:A:LEU:H	1:166:A:ASP:H	5	0.21
(1,13)	1:96:A:SER:H	1:157:A:VAL:H	20	0.21
(3,164)	1:90:A:GLN:H	1:177:A:VAL:H	4	0.2

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,109)	1:26:A:ILE:HD11	1:177:A:VAL:HG11	19	0.2
(3,109)	1:26:A:ILE:HD11	1:177:A:VAL:HG12	19	0.2
(3,109)	1:26:A:ILE:HD11	1:177:A:VAL:HG13	19	0.2
(3,109)	1:26:A:ILE:HD12	1:177:A:VAL:HG11	19	0.2
(3,109)	1:26:A:ILE:HD12	1:177:A:VAL:HG12	19	0.2
(3,109)	1:26:A:ILE:HD12	1:177:A:VAL:HG13	19	0.2
(3,109)	1:26:A:ILE:HD13	1:177:A:VAL:HG11	19	0.2
(3,109)	1:26:A:ILE:HD13	1:177:A:VAL:HG12	19	0.2
(3,109)	1:26:A:ILE:HD13	1:177:A:VAL:HG13	19	0.2
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	13	0.2
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	13	0.2
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	13	0.2
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	13	0.2
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	13	0.2
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	13	0.2
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	13	0.2
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	13	0.2
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	13	0.2
(3,46)	1:90:A:GLN:H	1:177:A:VAL:H	4	0.2
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG21	5	0.2
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG22	5	0.2
(1,242)	1:91:A:ILE:HD11	1:177:A:VAL:HG23	5	0.2
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG21	5	0.2
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG22	5	0.2
(1,242)	1:91:A:ILE:HD12	1:177:A:VAL:HG23	5	0.2
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG21	5	0.2
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG22	5	0.2
(1,242)	1:91:A:ILE:HD13	1:177:A:VAL:HG23	5	0.2
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	5	0.2
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	6	0.2
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	8	0.2
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	18	0.2
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	19	0.2
(1,203)	1:50:A:LEU:H	1:51:A:ALA:H	10	0.2
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	18	0.2
(1,179)	1:90:A:GLN:H	1:176:A:PHE:H	11	0.2
(1,159)	1:46:A:THR:H	1:49:A:ASP:H	7	0.2
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	13	0.2
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	1	0.2
(1,104)	1:84:A:ILE:H	1:91:A:ILE:H	1	0.2
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	6	0.2
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	17	0.2

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	20	0.2
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	5	0.2
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	19	0.2
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	7	0.2
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	6	0.19
(3,171)	1:136:A:MET:H	1:141:A:ILE:H	19	0.19
(3,164)	1:90:A:GLN:H	1:177:A:VAL:H	20	0.19
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	19	0.19
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	9	0.19
(3,151)	1:45:A:LEU:H	1:124:A:ASP:H	7	0.19
(3,128)	1:119:A:LEU:H	1:163:A:PHE:H	16	0.19
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	18	0.19
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	1	0.19
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	1	0.19
(3,80)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	1	0.19
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	1	0.19
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	1	0.19
(3,80)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	1	0.19
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	1	0.19
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	1	0.19
(3,80)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	1	0.19
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	6	0.19
(3,53)	1:136:A:MET:H	1:141:A:ILE:H	19	0.19
(3,46)	1:90:A:GLN:H	1:177:A:VAL:H	20	0.19
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	19	0.19
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	9	0.19
(3,33)	1:45:A:LEU:H	1:124:A:ASP:H	7	0.19
(3,10)	1:119:A:LEU:H	1:163:A:PHE:H	16	0.19
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	18	0.19
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	2	0.19
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	2	0.19
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	2	0.19
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	2	0.19
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	2	0.19
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	2	0.19
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	2	0.19
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	2	0.19
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	2	0.19
(1,149)	1:68:A:LYS:H	1:85:A:ARG:H	6	0.19
(1,145)	1:118:A:CYS:H	1:119:A:LEU:H	16	0.19
(1,129)	1:180:A:TYR:H	1:181:A:LEU:H	5	0.19
(1,126)	1:86:A:ASP:H	1:88:A:SER:H	15	0.19

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,118)	1:68:A:LYS:H	1:85:A:ARG:H	6	0.19
(1,104)	1:84:A:ILE:H	1:91:A:ILE:H	2	0.19
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	2	0.19
(1,55)	1:88:A:SER:H	1:89:A:HIS:H	11	0.19
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	19	0.19
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	2	0.19
(1,15)	1:180:A:TYR:H	1:181:A:LEU:H	5	0.19
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	6	0.19
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	16	0.18
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	14	0.18
(3,145)	1:42:A:ILE:H	1:125:A:ALA:H	1	0.18
(3,137)	1:42:A:ILE:H	1:125:A:ALA:H	1	0.18
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	6	0.18
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	16	0.18
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	16	0.18
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	14	0.18
(3,27)	1:42:A:ILE:H	1:125:A:ALA:H	1	0.18
(3,19)	1:42:A:ILE:H	1:125:A:ALA:H	1	0.18
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	6	0.18
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	16	0.18
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG11	7	0.18
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG12	7	0.18
(1,274)	1:26:A:ILE:HD11	1:177:A:VAL:HG13	7	0.18
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG11	7	0.18
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG12	7	0.18
(1,274)	1:26:A:ILE:HD12	1:177:A:VAL:HG13	7	0.18
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG11	7	0.18
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG12	7	0.18
(1,274)	1:26:A:ILE:HD13	1:177:A:VAL:HG13	7	0.18
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	10	0.18
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	10	0.18
(1,227)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	10	0.18
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	10	0.18
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	10	0.18
(1,227)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	10	0.18
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	10	0.18
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	10	0.18
(1,227)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	10	0.18
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	1	0.18
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	14	0.18
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	20	0.18
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	5	0.18

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	16	0.18
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	16	0.18
(1,91)	1:23:A:LYS:H	1:25:A:LEU:H	18	0.18
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	16	0.18
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	3	0.18
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	8	0.18
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	13	0.18
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	16	0.18
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	20	0.18
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	2	0.18
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	11	0.18
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	16	0.18
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	17	0.17
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	7	0.17
(3,120)	1:119:A:LEU:H	1:163:A:PHE:H	18	0.17
(3,108)	1:130:A:VAL:HG11	1:152:A:VAL:HG21	10	0.17
(3,108)	1:130:A:VAL:HG11	1:152:A:VAL:HG22	10	0.17
(3,108)	1:130:A:VAL:HG11	1:152:A:VAL:HG23	10	0.17
(3,108)	1:130:A:VAL:HG12	1:152:A:VAL:HG21	10	0.17
(3,108)	1:130:A:VAL:HG12	1:152:A:VAL:HG22	10	0.17
(3,108)	1:130:A:VAL:HG12	1:152:A:VAL:HG23	10	0.17
(3,108)	1:130:A:VAL:HG13	1:152:A:VAL:HG21	10	0.17
(3,108)	1:130:A:VAL:HG13	1:152:A:VAL:HG22	10	0.17
(3,108)	1:130:A:VAL:HG13	1:152:A:VAL:HG23	10	0.17
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	17	0.17
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	7	0.17
(3,2)	1:119:A:LEU:H	1:163:A:PHE:H	18	0.17
(1,246)	1:130:A:VAL:HG11	1:152:A:VAL:HG21	10	0.17
(1,246)	1:130:A:VAL:HG11	1:152:A:VAL:HG22	10	0.17
(1,246)	1:130:A:VAL:HG11	1:152:A:VAL:HG23	10	0.17
(1,246)	1:130:A:VAL:HG12	1:152:A:VAL:HG21	10	0.17
(1,246)	1:130:A:VAL:HG12	1:152:A:VAL:HG22	10	0.17
(1,246)	1:130:A:VAL:HG12	1:152:A:VAL:HG23	10	0.17
(1,246)	1:130:A:VAL:HG13	1:152:A:VAL:HG21	10	0.17
(1,246)	1:130:A:VAL:HG13	1:152:A:VAL:HG22	10	0.17
(1,246)	1:130:A:VAL:HG13	1:152:A:VAL:HG23	10	0.17
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	18	0.17
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	18	0.17
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	18	0.17
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	18	0.17
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	18	0.17
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	18	0.17

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	18	0.17
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	18	0.17
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	18	0.17
(1,210)	1:93:A:VAL:H	1:94:A:GLU:H	13	0.17
(1,208)	1:74:A:ASN:H	1:75:A:VAL:H	6	0.17
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	1	0.17
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	6	0.17
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	13	0.17
(1,130)	1:130:A:VAL:H	1:152:A:VAL:H	7	0.17
(1,126)	1:86:A:ASP:H	1:88:A:SER:H	6	0.17
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	20	0.17
(1,104)	1:84:A:ILE:H	1:91:A:ILE:H	10	0.17
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	18	0.17
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	7	0.17
(1,40)	1:124:A:ASP:H	1:125:A:ALA:H	15	0.17
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	14	0.17
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	4	0.17
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	9	0.17
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	9	0.16
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	9	0.16
(3,160)	1:84:A:ILE:H	1:91:A:ILE:H	20	0.16
(3,157)	1:44:A:SER:H	1:124:A:ASP:H	18	0.16
(3,134)	1:131:A:THR:H	1:135:A:ALA:H	5	0.16
(3,126)	1:126:A:ASP:H	1:156:A:ASN:H	12	0.16
(3,77)	1:91:A:ILE:HD11	1:177:A:VAL:HG21	16	0.16
(3,77)	1:91:A:ILE:HD11	1:177:A:VAL:HG22	16	0.16
(3,77)	1:91:A:ILE:HD11	1:177:A:VAL:HG23	16	0.16
(3,77)	1:91:A:ILE:HD12	1:177:A:VAL:HG21	16	0.16
(3,77)	1:91:A:ILE:HD12	1:177:A:VAL:HG22	16	0.16
(3,77)	1:91:A:ILE:HD12	1:177:A:VAL:HG23	16	0.16
(3,77)	1:91:A:ILE:HD13	1:177:A:VAL:HG21	16	0.16
(3,77)	1:91:A:ILE:HD13	1:177:A:VAL:HG22	16	0.16
(3,77)	1:91:A:ILE:HD13	1:177:A:VAL:HG23	16	0.16
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	9	0.16
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	9	0.16
(3,42)	1:84:A:ILE:H	1:91:A:ILE:H	20	0.16
(3,39)	1:44:A:SER:H	1:124:A:ASP:H	18	0.16
(3,16)	1:131:A:THR:H	1:135:A:ALA:H	5	0.16
(3,8)	1:126:A:ASP:H	1:156:A:ASN:H	12	0.16
(2,6)	1:18:A:LEU:HD21	1:119:A:LEU:HD11	8	0.16
(2,6)	1:18:A:LEU:HD21	1:119:A:LEU:HD12	8	0.16
(2,6)	1:18:A:LEU:HD21	1:119:A:LEU:HD13	8	0.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,6)	1:18:A:LEU:HD22	1:119:A:LEU:HD11	8	0.16
(2,6)	1:18:A:LEU:HD22	1:119:A:LEU:HD12	8	0.16
(2,6)	1:18:A:LEU:HD22	1:119:A:LEU:HD13	8	0.16
(2,6)	1:18:A:LEU:HD23	1:119:A:LEU:HD11	8	0.16
(2,6)	1:18:A:LEU:HD23	1:119:A:LEU:HD12	8	0.16
(2,6)	1:18:A:LEU:HD23	1:119:A:LEU:HD13	8	0.16
(1,282)	1:41:A:VAL:HG21	1:155:A:LEU:HD11	4	0.16
(1,282)	1:41:A:VAL:HG21	1:155:A:LEU:HD12	4	0.16
(1,282)	1:41:A:VAL:HG21	1:155:A:LEU:HD13	4	0.16
(1,282)	1:41:A:VAL:HG22	1:155:A:LEU:HD11	4	0.16
(1,282)	1:41:A:VAL:HG22	1:155:A:LEU:HD12	4	0.16
(1,282)	1:41:A:VAL:HG22	1:155:A:LEU:HD13	4	0.16
(1,282)	1:41:A:VAL:HG23	1:155:A:LEU:HD11	4	0.16
(1,282)	1:41:A:VAL:HG23	1:155:A:LEU:HD12	4	0.16
(1,282)	1:41:A:VAL:HG23	1:155:A:LEU:HD13	4	0.16
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD11	20	0.16
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD12	20	0.16
(1,278)	1:41:A:VAL:HG21	1:155:A:LEU:HD13	20	0.16
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD11	20	0.16
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD12	20	0.16
(1,278)	1:41:A:VAL:HG22	1:155:A:LEU:HD13	20	0.16
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD11	20	0.16
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD12	20	0.16
(1,278)	1:41:A:VAL:HG23	1:155:A:LEU:HD13	20	0.16
(1,203)	1:50:A:LEU:H	1:51:A:ALA:H	2	0.16
(1,186)	1:76:A:CYS:H	1:78:A:TYR:H	18	0.16
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	7	0.16
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	3	0.16
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	5	0.16
(1,110)	1:45:A:LEU:H	1:124:A:ASP:H	11	0.16
(1,107)	1:67:A:ALA:H	1:120:A:MET:H	8	0.16
(1,94)	1:121:A:ILE:H	1:163:A:PHE:H	9	0.16
(1,88)	1:42:A:ILE:H	1:125:A:ALA:H	16	0.16
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	9	0.16
(1,63)	1:86:A:ASP:H	1:88:A:SER:H	14	0.16
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	2	0.16
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	10	0.16
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	12	0.16
(1,41)	1:42:A:ILE:H	1:125:A:ALA:H	16	0.16
(1,20)	1:76:A:CYS:H	1:79:A:LYS:H	20	0.16
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	7	0.16
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	10	0.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	18	0.16
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	14	0.15
(3,166)	1:90:A:GLN:H	1:178:A:TYR:H	16	0.15
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	5	0.15
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	3	0.15
(3,132)	1:41:A:VAL:H	1:127:A:LEU:H	18	0.15
(3,120)	1:119:A:LEU:H	1:163:A:PHE:H	11	0.15
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	14	0.15
(3,48)	1:90:A:GLN:H	1:178:A:TYR:H	16	0.15
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	5	0.15
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	3	0.15
(3,14)	1:41:A:VAL:H	1:127:A:LEU:H	18	0.15
(3,2)	1:119:A:LEU:H	1:163:A:PHE:H	11	0.15
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD11	15	0.15
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD12	15	0.15
(1,267)	1:18:A:LEU:HD11	1:119:A:LEU:HD13	15	0.15
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD11	15	0.15
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD12	15	0.15
(1,267)	1:18:A:LEU:HD12	1:119:A:LEU:HD13	15	0.15
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD11	15	0.15
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD12	15	0.15
(1,267)	1:18:A:LEU:HD13	1:119:A:LEU:HD13	15	0.15
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD11	8	0.15
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD12	8	0.15
(1,220)	1:69:A:ILE:HD11	1:111:A:ILE:HD13	8	0.15
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD11	8	0.15
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD12	8	0.15
(1,220)	1:69:A:ILE:HD12	1:111:A:ILE:HD13	8	0.15
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD11	8	0.15
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD12	8	0.15
(1,220)	1:69:A:ILE:HD13	1:111:A:ILE:HD13	8	0.15
(1,211)	1:131:A:THR:H	1:134:A:ARG:H	6	0.15
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	4	0.15
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	16	0.15
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	4	0.15
(1,182)	1:108:A:ASN:H	1:109:A:CYS:H	12	0.15
(1,182)	1:108:A:ASN:H	1:109:A:CYS:H	19	0.15
(1,178)	1:90:A:GLN:H	1:177:A:VAL:H	9	0.15
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	3	0.15
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	8	0.15
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	6	0.15
(1,110)	1:45:A:LEU:H	1:124:A:ASP:H	19	0.15

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,75)	1:90:A:GLN:H	1:177:A:VAL:H	7	0.15
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	18	0.15
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	8	0.15
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	17	0.15
(1,22)	1:119:A:LEU:H	1:166:A:ASP:H	5	0.15
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	19	0.15
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	14	0.14
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	3	0.14
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	11	0.14
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	13	0.14
(3,165)	1:90:A:GLN:H	1:176:A:PHE:H	3	0.14
(3,160)	1:84:A:ILE:H	1:91:A:ILE:H	1	0.14
(3,134)	1:131:A:THR:H	1:135:A:ALA:H	3	0.14
(3,134)	1:131:A:THR:H	1:135:A:ALA:H	11	0.14
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	13	0.14
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	13	0.14
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	13	0.14
(3,63)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	13	0.14
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	13	0.14
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	13	0.14
(3,63)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	13	0.14
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	13	0.14
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	13	0.14
(3,63)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	13	0.14
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	14	0.14
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	3	0.14
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	11	0.14
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	13	0.14
(3,47)	1:90:A:GLN:H	1:176:A:PHE:H	3	0.14
(3,42)	1:84:A:ILE:H	1:91:A:ILE:H	1	0.14
(3,16)	1:131:A:THR:H	1:135:A:ALA:H	3	0.14
(3,16)	1:131:A:THR:H	1:135:A:ALA:H	11	0.14
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	13	0.14
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	17	0.14
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	17	0.14
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	17	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	17	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	17	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	17	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	17	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	17	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	17	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	18	0.14
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	18	0.14
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	18	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	18	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	18	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	18	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	18	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	18	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	18	0.14
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	19	0.14
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	19	0.14
(1,254)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	19	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	19	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	19	0.14
(1,254)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	19	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	19	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	19	0.14
(1,254)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	19	0.14
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	9	0.14
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	19	0.14
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	20	0.14
(1,159)	1:46:A:THR:H	1:49:A:ASP:H	8	0.14
(1,156)	1:84:A:ILE:H	1:91:A:ILE:H	18	0.14
(1,149)	1:68:A:LYS:H	1:85:A:ARG:H	3	0.14
(1,118)	1:68:A:LYS:H	1:85:A:ARG:H	3	0.14
(1,104)	1:84:A:ILE:H	1:91:A:ILE:H	14	0.14
(1,62)	1:86:A:ASP:H	1:89:A:HIS:H	16	0.14
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	4	0.14
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	11	0.14
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	15	0.14
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	20	0.14
(1,34)	1:131:A:THR:H	1:135:A:ALA:H	7	0.14
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	1	0.13
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	2	0.13
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	16	0.13
(3,169)	1:46:A:THR:H	1:50:A:LEU:H	18	0.13
(3,160)	1:84:A:ILE:H	1:91:A:ILE:H	2	0.13
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	4	0.13
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	11	0.13
(3,138)	1:128:A:VAL:H	1:154:A:VAL:H	18	0.13
(3,127)	1:96:A:SER:H	1:157:A:VAL:H	14	0.13
(3,112)	1:41:A:VAL:HG21	1:155:A:LEU:HD11	4	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,112)	1:41:A:VAL:HG21	1:155:A:LEU:HD12	4	0.13
(3,112)	1:41:A:VAL:HG21	1:155:A:LEU:HD13	4	0.13
(3,112)	1:41:A:VAL:HG22	1:155:A:LEU:HD11	4	0.13
(3,112)	1:41:A:VAL:HG22	1:155:A:LEU:HD12	4	0.13
(3,112)	1:41:A:VAL:HG22	1:155:A:LEU:HD13	4	0.13
(3,112)	1:41:A:VAL:HG23	1:155:A:LEU:HD11	4	0.13
(3,112)	1:41:A:VAL:HG23	1:155:A:LEU:HD12	4	0.13
(3,112)	1:41:A:VAL:HG23	1:155:A:LEU:HD13	4	0.13
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD21	9	0.13
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD22	9	0.13
(3,91)	1:146:A:ILE:HD11	1:171:A:LEU:HD23	9	0.13
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD21	9	0.13
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD22	9	0.13
(3,91)	1:146:A:ILE:HD12	1:171:A:LEU:HD23	9	0.13
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD21	9	0.13
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD22	9	0.13
(3,91)	1:146:A:ILE:HD13	1:171:A:LEU:HD23	9	0.13
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	1	0.13
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	2	0.13
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	16	0.13
(3,51)	1:46:A:THR:H	1:50:A:LEU:H	18	0.13
(3,42)	1:84:A:ILE:H	1:91:A:ILE:H	2	0.13
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	4	0.13
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	11	0.13
(3,20)	1:128:A:VAL:H	1:154:A:VAL:H	18	0.13
(3,9)	1:96:A:SER:H	1:157:A:VAL:H	14	0.13
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	1	0.13
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	1	0.13
(1,229)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	1	0.13
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	1	0.13
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	1	0.13
(1,229)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	1	0.13
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	1	0.13
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	1	0.13
(1,229)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	1	0.13
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	10	0.13
(1,205)	1:46:A:THR:H	1:50:A:LEU:H	20	0.13
(1,183)	1:137:A:SER:H	1:140:A:LYS:H	14	0.13
(1,182)	1:108:A:ASN:H	1:109:A:CYS:H	6	0.13
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	2	0.13
(1,75)	1:90:A:GLN:H	1:177:A:VAL:H	4	0.13
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	3	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	14	0.13
(1,33)	1:94:A:GLU:H	1:155:A:LEU:H	1	0.13
(1,33)	1:94:A:GLU:H	1:155:A:LEU:H	16	0.13
(1,23)	1:41:A:VAL:H	1:127:A:LEU:H	13	0.13
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	9	0.13
(3,213)	1:141:A:ILE:CB	1:171:A:LEU:CB	14	0.12
(3,213)	1:141:A:ILE:CB	1:171:A:LEU:CD1	14	0.12
(3,213)	1:141:A:ILE:CB	1:171:A:LEU:CD2	14	0.12
(3,213)	1:141:A:ILE:CB	1:171:A:LEU:CG	14	0.12
(3,213)	1:141:A:ILE:CD1	1:171:A:LEU:CB	14	0.12
(3,213)	1:141:A:ILE:CD1	1:171:A:LEU:CD1	14	0.12
(3,213)	1:141:A:ILE:CD1	1:171:A:LEU:CD2	14	0.12
(3,213)	1:141:A:ILE:CD1	1:171:A:LEU:CG	14	0.12
(3,213)	1:141:A:ILE:CG1	1:171:A:LEU:CB	14	0.12
(3,213)	1:141:A:ILE:CG1	1:171:A:LEU:CD1	14	0.12
(3,213)	1:141:A:ILE:CG1	1:171:A:LEU:CD2	14	0.12
(3,213)	1:141:A:ILE:CG1	1:171:A:LEU:CG	14	0.12
(3,213)	1:141:A:ILE:CG2	1:171:A:LEU:CB	14	0.12
(3,213)	1:141:A:ILE:CG2	1:171:A:LEU:CD1	14	0.12
(3,213)	1:141:A:ILE:CG2	1:171:A:LEU:CD2	14	0.12
(3,213)	1:141:A:ILE:CG2	1:171:A:LEU:CG	14	0.12
(3,173)	1:41:A:VAL:H	1:127:A:LEU:H	11	0.12
(3,152)	1:123:A:GLY:H	1:159:A:GLN:H	13	0.12
(3,150)	1:67:A:ALA:H	1:120:A:MET:H	1	0.12
(3,147)	1:121:A:ILE:H	1:163:A:PHE:H	5	0.12
(3,126)	1:126:A:ASP:H	1:156:A:ASN:H	8	0.12
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD21	1	0.12
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD22	1	0.12
(3,90)	1:141:A:ILE:HD11	1:171:A:LEU:HD23	1	0.12
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD21	1	0.12
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD22	1	0.12
(3,90)	1:141:A:ILE:HD12	1:171:A:LEU:HD23	1	0.12
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD21	1	0.12
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD22	1	0.12
(3,90)	1:141:A:ILE:HD13	1:171:A:LEU:HD23	1	0.12
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD11	9	0.12
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD12	9	0.12
(3,57)	1:69:A:ILE:HD11	1:111:A:ILE:HD13	9	0.12
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD11	9	0.12
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD12	9	0.12
(3,57)	1:69:A:ILE:HD12	1:111:A:ILE:HD13	9	0.12
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD11	9	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD12	9	0.12
(3,57)	1:69:A:ILE:HD13	1:111:A:ILE:HD13	9	0.12
(3,55)	1:41:A:VAL:H	1:127:A:LEU:H	11	0.12
(3,34)	1:123:A:GLY:H	1:159:A:GLN:H	13	0.12
(3,32)	1:67:A:ALA:H	1:120:A:MET:H	1	0.12
(3,29)	1:121:A:ILE:H	1:163:A:PHE:H	5	0.12
(3,8)	1:126:A:ASP:H	1:156:A:ASN:H	8	0.12
(1,196)	1:123:A:GLY:H	1:159:A:GLN:H	6	0.12
(1,159)	1:46:A:THR:H	1:49:A:ASP:H	17	0.12
(1,134)	1:178:A:TYR:H	1:180:A:TYR:H	11	0.12
(1,112)	1:123:A:GLY:H	1:159:A:GLN:H	1	0.12
(1,75)	1:90:A:GLN:H	1:177:A:VAL:H	20	0.12
(1,56)	1:86:A:ASP:H	1:89:A:HIS:H	20	0.12
(1,50)	1:71:A:LYS:H	1:72:A:PHE:H	6	0.12
(1,40)	1:124:A:ASP:H	1:125:A:ALA:H	4	0.12
(1,33)	1:94:A:GLU:H	1:155:A:LEU:H	8	0.12
(1,12)	1:126:A:ASP:H	1:156:A:ASN:H	17	0.12
(3,160)	1:84:A:ILE:H	1:91:A:ILE:H	10	0.11
(3,159)	1:41:A:VAL:H	1:125:A:ALA:H	6	0.11
(3,134)	1:131:A:THR:H	1:135:A:ALA:H	4	0.11
(3,124)	1:75:A:VAL:H	1:81:A:TYR:H	14	0.11
(3,42)	1:84:A:ILE:H	1:91:A:ILE:H	10	0.11
(3,41)	1:41:A:VAL:H	1:125:A:ALA:H	6	0.11
(3,16)	1:131:A:THR:H	1:135:A:ALA:H	4	0.11
(3,6)	1:75:A:VAL:H	1:81:A:TYR:H	14	0.11
(1,244)	1:69:A:ILE:HD11	1:111:A:ILE:HD11	13	0.11
(1,244)	1:69:A:ILE:HD11	1:111:A:ILE:HD12	13	0.11
(1,244)	1:69:A:ILE:HD11	1:111:A:ILE:HD13	13	0.11
(1,244)	1:69:A:ILE:HD12	1:111:A:ILE:HD11	13	0.11
(1,244)	1:69:A:ILE:HD12	1:111:A:ILE:HD12	13	0.11
(1,244)	1:69:A:ILE:HD12	1:111:A:ILE:HD13	13	0.11
(1,244)	1:69:A:ILE:HD13	1:111:A:ILE:HD11	13	0.11
(1,244)	1:69:A:ILE:HD13	1:111:A:ILE:HD12	13	0.11
(1,244)	1:69:A:ILE:HD13	1:111:A:ILE:HD13	13	0.11
(1,199)	1:16:A:VAL:H	1:18:A:LEU:H	16	0.11
(1,182)	1:108:A:ASN:H	1:109:A:CYS:H	15	0.11
(1,178)	1:90:A:GLN:H	1:177:A:VAL:H	15	0.11
(1,175)	1:137:A:SER:H	1:139:A:PHE:H	20	0.11
(1,133)	1:107:A:HIS:H	1:109:A:CYS:H	20	0.11
(1,104)	1:84:A:ILE:H	1:91:A:ILE:H	13	0.11
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	4	0.11
(1,99)	1:76:A:CYS:H	1:79:A:LYS:H	10	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	10	0.11
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	12	0.11
(1,40)	1:124:A:ASP:H	1:125:A:ALA:H	11	0.11
(1,40)	1:124:A:ASP:H	1:125:A:ALA:H	13	0.11
(1,29)	1:134:A:ARG:H	1:136:A:MET:H	3	0.11
(3,168)	1:123:A:GLY:H	1:159:A:GLN:H	11	0.1
(3,126)	1:126:A:ASP:H	1:156:A:ASN:H	5	0.1
(3,126)	1:126:A:ASP:H	1:156:A:ASN:H	15	0.1
(3,50)	1:123:A:GLY:H	1:159:A:GLN:H	11	0.1
(3,8)	1:126:A:ASP:H	1:156:A:ASN:H	5	0.1
(3,8)	1:126:A:ASP:H	1:156:A:ASN:H	15	0.1
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD11	20	0.1
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD12	20	0.1
(1,245)	1:92:A:LEU:HD11	1:141:A:ILE:HD13	20	0.1
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD11	20	0.1
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD12	20	0.1
(1,245)	1:92:A:LEU:HD12	1:141:A:ILE:HD13	20	0.1
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD11	20	0.1
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD12	20	0.1
(1,245)	1:92:A:LEU:HD13	1:141:A:ILE:HD13	20	0.1
(1,219)	1:94:A:GLU:H	1:155:A:LEU:H	16	0.1
(1,218)	1:41:A:VAL:H	1:127:A:LEU:H	2	0.1
(1,159)	1:46:A:THR:H	1:49:A:ASP:H	2	0.1
(1,159)	1:46:A:THR:H	1:49:A:ASP:H	3	0.1
(1,123)	1:177:A:VAL:H	1:179:A:LYS:H	11	0.1
(1,107)	1:67:A:ALA:H	1:120:A:MET:H	3	0.1
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	1	0.1
(1,71)	1:130:A:VAL:H	1:152:A:VAL:H	9	0.1
(1,3)	1:119:A:LEU:H	1:163:A:PHE:H	20	0.1

10 Dihedral-angle violation analysis ⓘ

No dihedral-angle restraints found