



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 07:31 AM UTC

PDB ID : 2M83 / pdb_00002m83
BMRB ID : 19225
Title : Solution structure of the carbohydrate binding module of the muscle glycogen-targeting subunit of Protein Phosphatase-1
Authors : Koveal, D.; Page, R.; Peti, W.
Deposited on : 2013-05-03

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

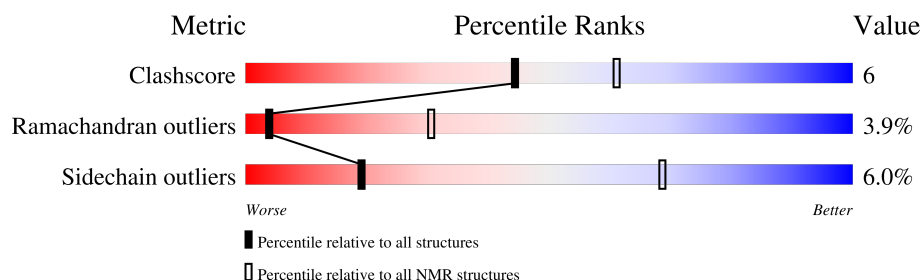
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 82%.

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	139	

2 Ensemble composition and analysis

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

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

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:106-A:237 (132)	0.81	13

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

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

The models can be grouped into 3 clusters and 6 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2211 atoms, of which 1089 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Protein phosphatase 1 regulatory subunit 3A.

Mol	Chain	Residues	Atoms						Trace
1	A	139	Total	C	H	N	O	S	0
			2211	713	1089	178	223	8	

There are 3 discrepancies between the modelled and reference sequences:

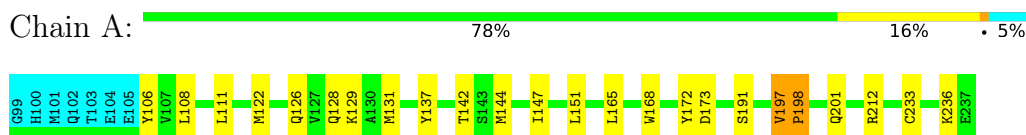
Chain	Residue	Modelled	Actual	Comment	Reference
A	99	GLY	-	expression tag	UNP Q00756
A	100	HIS	-	expression tag	UNP Q00756
A	101	MET	-	expression tag	UNP Q00756

4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

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

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A

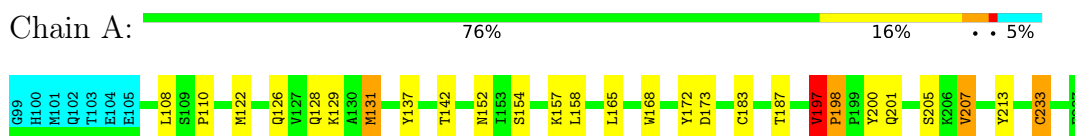


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

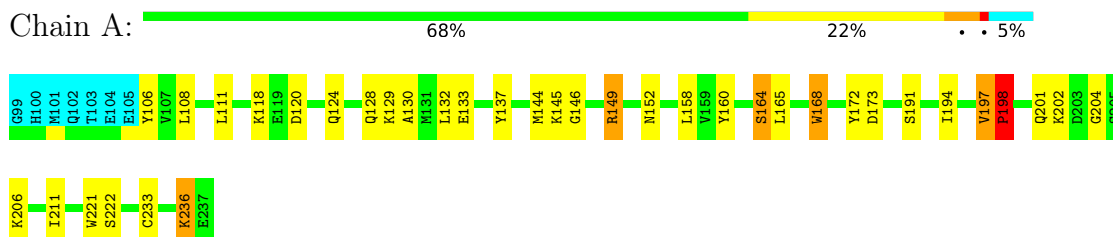
4.2.1 Score per residue for model 1

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



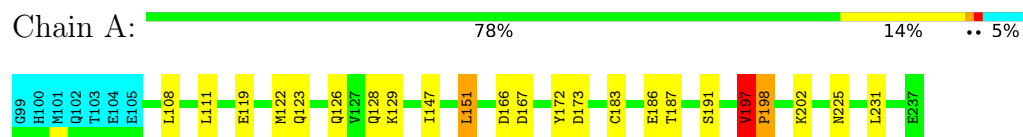
4.2.2 Score per residue for model 2

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



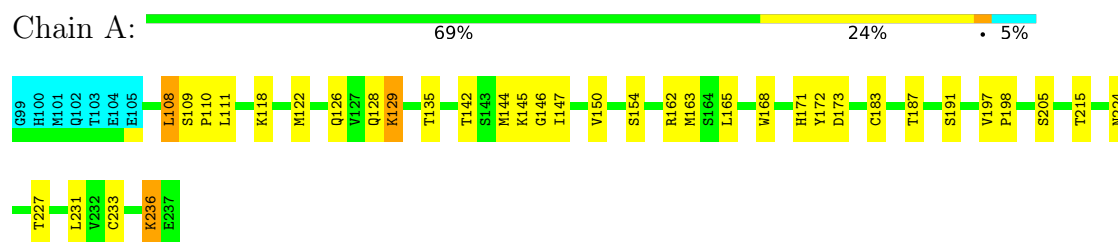
4.2.3 Score per residue for model 3

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



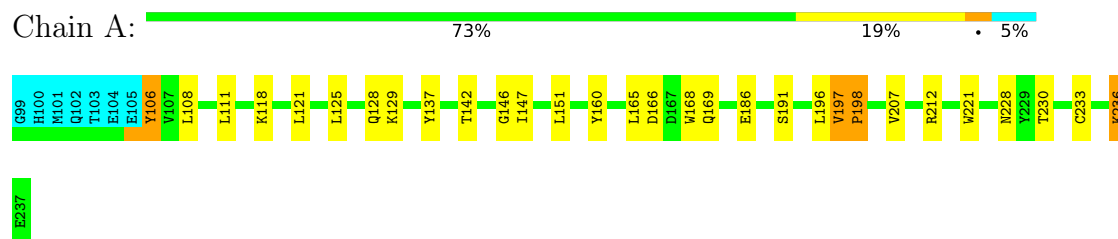
4.2.4 Score per residue for model 4

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



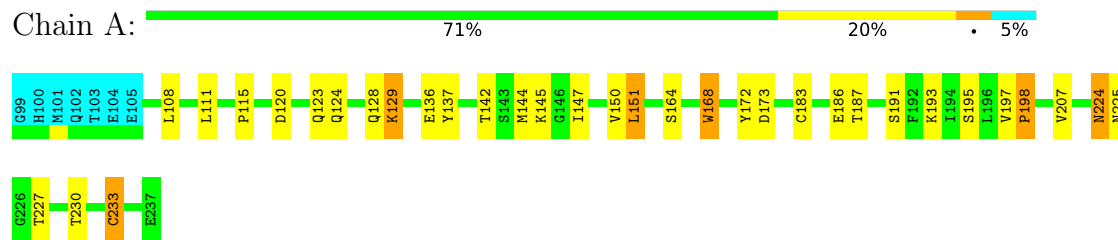
4.2.5 Score per residue for model 5

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



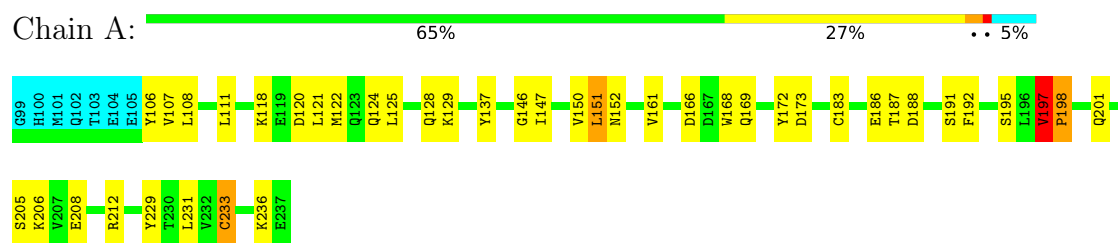
4.2.6 Score per residue for model 6

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



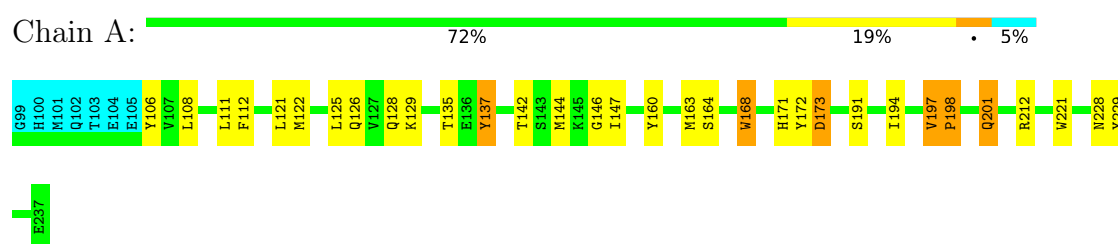
4.2.7 Score per residue for model 7

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



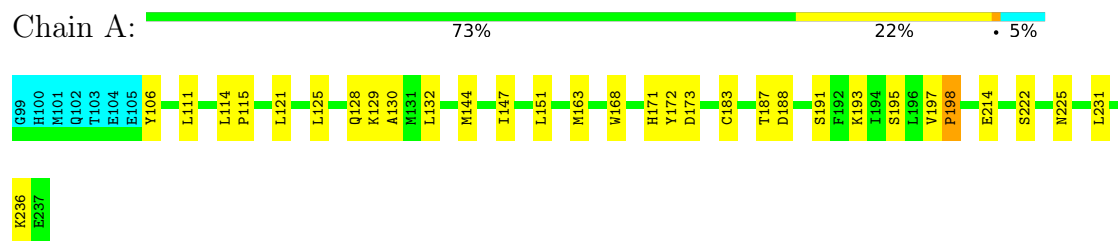
4.2.8 Score per residue for model 8

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



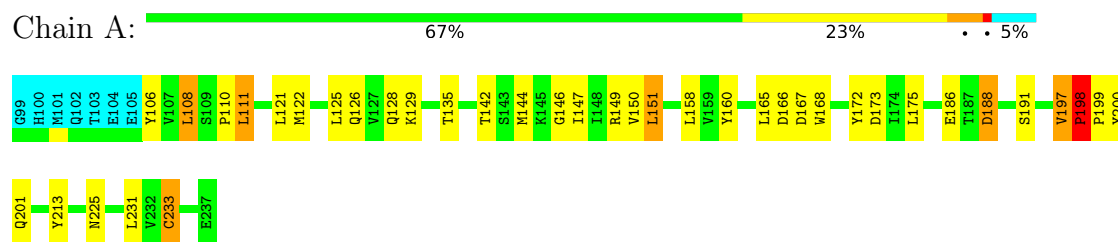
4.2.9 Score per residue for model 9

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



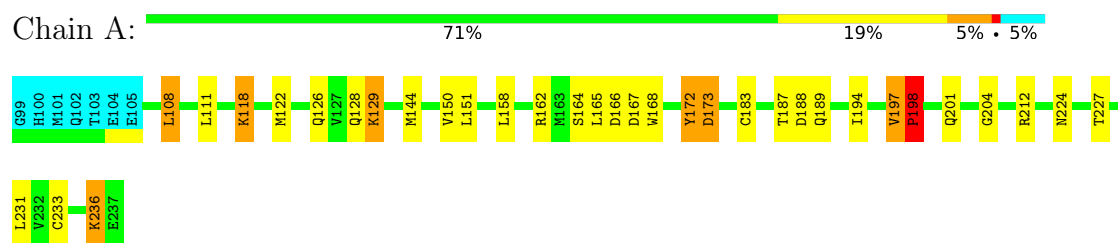
4.2.10 Score per residue for model 10

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



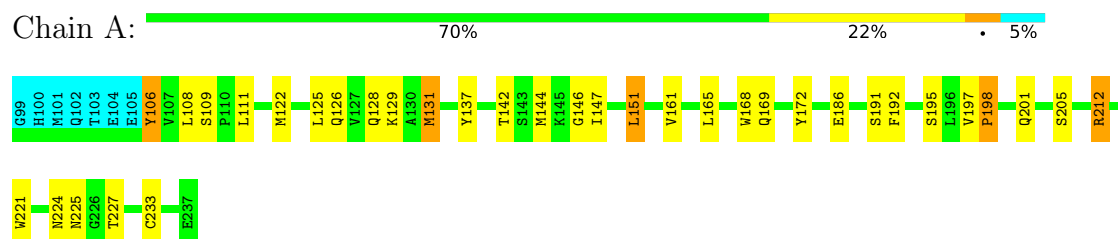
4.2.11 Score per residue for model 11

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



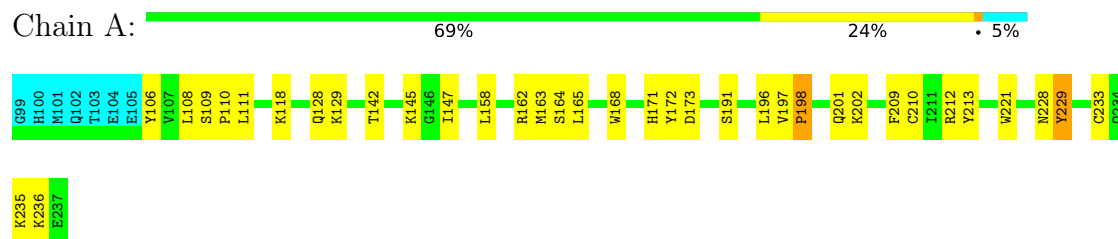
4.2.12 Score per residue for model 12

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



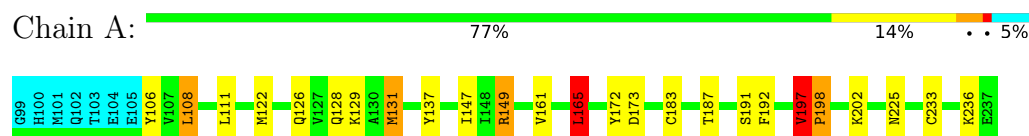
4.2.13 Score per residue for model 13 (medoid)

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



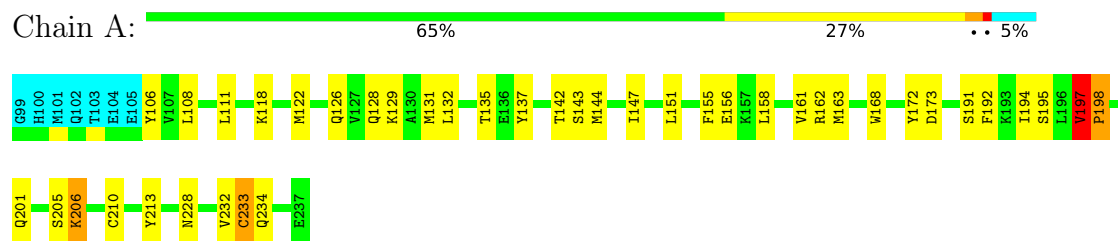
4.2.14 Score per residue for model 14

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



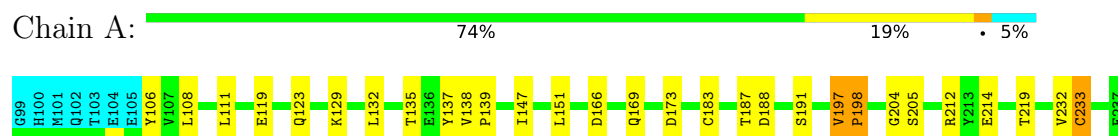
4.2.15 Score per residue for model 15

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



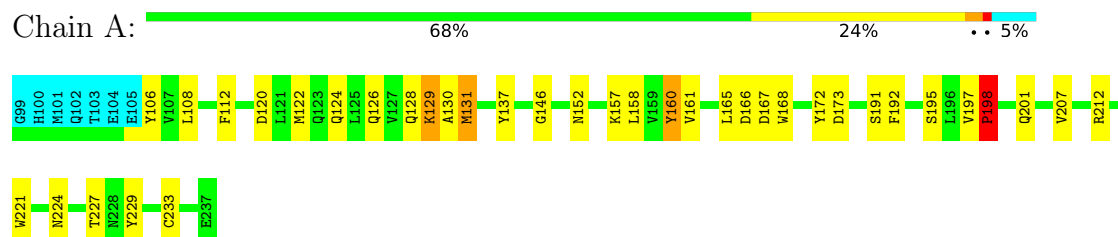
4.2.16 Score per residue for model 16

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



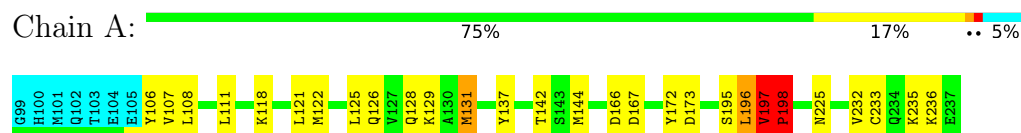
4.2.17 Score per residue for model 17

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



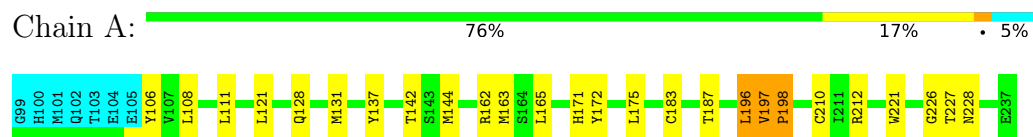
4.2.18 Score per residue for model 18

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



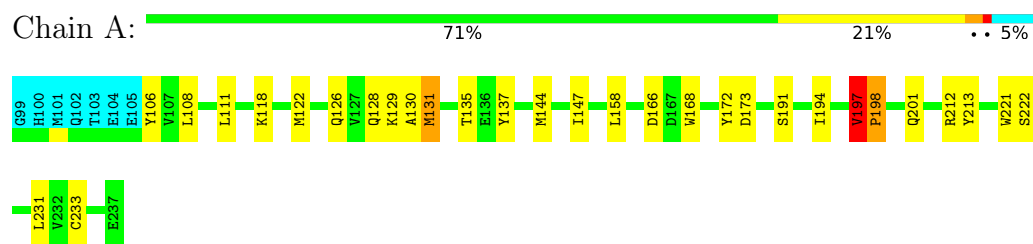
4.2.19 Score per residue for model 19

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



4.2.20 Score per residue for model 20

- Molecule 1: Protein phosphatase 1 regulatory subunit 3A



5 Refinement protocol and experimental data overview

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

Of the 200 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
CNS	refinement	
CYANA 2.1	structure solution	

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	1524
Number of shifts mapped to atoms	1524
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	82%

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	1.23±0.04	1±1/1090 (0.1± 0.1%)	1.02±0.03	4±1/1476 (0.2± 0.1%)
All	All	1.24	27/21800 (0.1%)	1.02	70/29520 (0.2%)

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

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	0.5±0.5
All	All	0	9

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	106	TYR	N-CA	-7.44	1.36	1.46	5	1
1	A	109	SER	N-CA	-7.02	1.38	1.46	13	3
1	A	152	ASN	C-N	-6.21	1.28	1.33	17	4
1	A	164	SER	N-CA	-6.21	1.39	1.46	11	3
1	A	149	ARG	C-N	-6.01	1.25	1.33	10	1
1	A	160	TYR	C-N	-6.00	1.25	1.33	17	1
1	A	214	GLU	C-N	-5.78	1.28	1.33	9	1
1	A	149	ARG	CA-C	-5.66	1.45	1.52	10	1
1	A	227	THR	C-N	-5.52	1.26	1.33	19	1
1	A	157	LYS	N-CA	-5.47	1.39	1.46	17	1
1	A	145	LYS	C-N	-5.43	1.29	1.33	2	2
1	A	197	VAL	C-N	5.41	1.39	1.33	16	2
1	A	227	THR	N-CA	-5.38	1.39	1.45	19	1
1	A	158	LEU	C-N	-5.32	1.28	1.33	17	1
1	A	226	GLY	C-N	-5.16	1.26	1.33	19	1
1	A	207	VAL	N-CA	-5.13	1.40	1.46	1	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	129	LYS	N-CA	5.03	1.52	1.46	17	1

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	198	PRO	N-CA-C	8.68	121.29	110.70	10	20
1	A	197	VAL	CA-C-N	7.79	128.41	120.38	10	15
1	A	197	VAL	C-N-CA	7.79	128.41	120.38	10	15
1	A	165	LEU	N-CA-C	-7.34	105.19	112.97	12	8
1	A	173	ASP	CA-CB-CG	6.70	119.30	112.60	11	1
1	A	166	ASP	CA-CB-CG	5.84	118.44	112.60	20	1
1	A	169	GLN	N-CA-C	-5.74	106.40	113.41	12	1
1	A	232	VAL	CB-CA-C	-5.49	104.28	111.25	18	1
1	A	145	LYS	CA-C-N	-5.35	118.49	122.18	4	1
1	A	145	LYS	C-N-CA	-5.35	118.49	122.18	4	1
1	A	138	VAL	CA-C-N	5.24	126.39	119.84	16	1
1	A	138	VAL	C-N-CA	5.24	126.39	119.84	16	1
1	A	204	GLY	CA-C-N	5.20	127.51	120.38	16	1
1	A	204	GLY	C-N-CA	5.20	127.51	120.38	16	1
1	A	228	ASN	N-CA-C	-5.09	105.30	112.12	5	1
1	A	188	ASP	CA-CB-CG	5.04	117.64	112.60	10	1

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	162	ARG	Sidechain	4
1	A	212	ARG	Sidechain	4
1	A	149	ARG	Sidechain	1

6.2 Too-close contacts ⓘ

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	1066	1040	1039	13±4
All	All	21320	20800	20780	268

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 6.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:197:VAL:HB	1:A:198:PRO:HD2	0.72	1.61	16	11
1:A:210:CYS:SG	1:A:228:ASN:HB3	0.65	2.32	13	3
1:A:198:PRO:HA	1:A:201:GLN:O	0.63	1.92	13	8
1:A:197:VAL:HB	1:A:198:PRO:CD	0.63	2.23	13	5
1:A:130:ALA:O	1:A:222:SER:HB3	0.62	1.93	20	2
1:A:106:TYR:HB3	1:A:137:TYR:CD2	0.60	2.31	7	3
1:A:224:ASN:HB3	1:A:227:THR:O	0.60	1.97	6	5
1:A:106:TYR:HB2	1:A:137:TYR:CD2	0.59	2.31	5	2
1:A:142:THR:HG23	1:A:200:TYR:CD1	0.59	2.33	10	1
1:A:147:ILE:HD13	1:A:191:SER:HB3	0.58	1.75	3	14
1:A:108:LEU:HD12	1:A:144:MET:SD	0.58	2.39	4	8
1:A:122:MET:O	1:A:126:GLN:HG2	0.57	2.00	15	12
1:A:183:CYS:HA	1:A:187:THR:O	0.56	1.99	3	10
1:A:160:TYR:CE2	1:A:173:ASP:HB2	0.56	2.34	8	1
1:A:211:ILE:O	1:A:221:TRP:HA	0.55	2.00	2	1
1:A:205:SER:HB3	1:A:233:CYS:SG	0.55	2.41	7	2
1:A:143:SER:HA	1:A:195:SER:O	0.55	2.02	15	1
1:A:205:SER:HB2	1:A:233:CYS:SG	0.54	2.43	15	4
1:A:108:LEU:HG	1:A:137:TYR:CE1	0.53	2.37	12	7
1:A:162:ARG:HG2	1:A:173:ASP:CB	0.53	2.34	11	1
1:A:212:ARG:HD2	1:A:221:TRP:CZ2	0.53	2.39	8	6
1:A:166:ASP:OD2	1:A:169:GLN:HG2	0.52	2.04	7	3
1:A:151:LEU:HD13	1:A:186:GLU:O	0.52	2.05	12	6
1:A:163:MET:O	1:A:171:HIS:HA	0.52	2.05	19	5
1:A:208:GLU:HA	1:A:229:TYR:O	0.52	2.05	7	1
1:A:144:MET:HB3	1:A:194:ILE:CG2	0.51	2.36	20	4
1:A:135:THR:HB	1:A:144:MET:SD	0.51	2.44	15	5
1:A:121:LEU:O	1:A:125:LEU:HB2	0.51	2.05	18	5
1:A:131:MET:HE3	1:A:131:MET:O	0.51	2.06	1	5
1:A:146:GLY:O	1:A:191:SER:HA	0.51	2.06	5	8
1:A:158:LEU:O	1:A:213:TYR:HA	0.51	2.05	1	5
1:A:207:VAL:O	1:A:230:THR:HA	0.49	2.08	5	2
1:A:108:LEU:HD22	1:A:231:LEU:HB3	0.49	1.83	7	2
1:A:183:CYS:HB3	1:A:188:ASP:OD1	0.48	2.08	16	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:111:LEU:HD11	1:A:232:VAL:HG13	0.48	1.84	16	2
1:A:108:LEU:HD21	1:A:233:CYS:HA	0.48	1.84	12	4
1:A:120:ASP:O	1:A:124:GLN:HG2	0.48	2.09	17	4
1:A:197:VAL:CB	1:A:198:PRO:HD2	0.48	2.39	1	13
1:A:161:VAL:HG11	1:A:192:PHE:CD2	0.47	2.45	12	4
1:A:198:PRO:HD3	1:A:233:CYS:SG	0.47	2.50	6	1
1:A:128:GLN:OE1	1:A:131:MET:HG2	0.47	2.09	17	1
1:A:158:LEU:HD12	1:A:160:TYR:CE2	0.47	2.45	10	1
1:A:154:SER:O	1:A:157:LYS:HE2	0.47	2.10	1	1
1:A:132:LEU:HD11	1:A:135:THR:HG23	0.46	1.87	15	2
1:A:107:VAL:HG22	1:A:236:LYS:HB2	0.46	1.88	7	2
1:A:144:MET:O	1:A:193:LYS:HA	0.46	2.11	9	2
1:A:106:TYR:CE1	1:A:235:LYS:HE2	0.46	2.46	13	1
1:A:108:LEU:HG	1:A:137:TYR:OH	0.46	2.11	14	2
1:A:198:PRO:HG3	1:A:233:CYS:C	0.45	2.36	13	1
1:A:112:PHE:CD1	1:A:229:TYR:HA	0.45	2.47	17	1
1:A:166:ASP:CG	1:A:167:ASP:H	0.45	2.20	11	5
1:A:111:LEU:H	1:A:111:LEU:HD23	0.45	1.71	18	1
1:A:198:PRO:HD3	1:A:233:CYS:HB3	0.45	1.88	17	1
1:A:197:VAL:C	1:A:199:PRO:HD2	0.45	2.37	10	1
1:A:163:MET:SD	1:A:195:SER:HB2	0.45	2.52	15	1
1:A:142:THR:HA	1:A:200:TYR:CE2	0.44	2.47	1	1
1:A:155:PHE:CD2	1:A:156:GLU:HG2	0.44	2.48	15	1
1:A:111:LEU:N	1:A:111:LEU:HD12	0.44	2.27	19	2
1:A:118:LYS:O	1:A:122:MET:HG3	0.43	2.13	11	1
1:A:196:LEU:HD12	1:A:196:LEU:H	0.43	1.73	18	1
1:A:164:SER:OG	1:A:168:TRP:HA	0.43	2.13	8	3
1:A:172:TYR:C	1:A:172:TYR:CD1	0.43	2.96	11	1
1:A:133:GLU:OE1	1:A:149:ARG:HG3	0.43	2.13	2	1
1:A:150:VAL:CG2	1:A:188:ASP:HB2	0.43	2.43	11	3
1:A:161:VAL:HG11	1:A:192:PHE:CE2	0.43	2.48	17	3
1:A:121:LEU:HG	1:A:131:MET:SD	0.43	2.54	19	1
1:A:151:LEU:HG	1:A:187:THR:HG22	0.43	1.90	11	2
1:A:162:ARG:HG2	1:A:173:ASP:HB3	0.43	1.91	11	1
1:A:107:VAL:O	1:A:235:LYS:HA	0.43	2.14	18	1
1:A:108:LEU:O	1:A:110:PRO:HD3	0.42	2.14	10	4
1:A:224:ASN:HB2	1:A:227:THR:O	0.42	2.15	4	1
1:A:108:LEU:CD2	1:A:233:CYS:HA	0.42	2.44	12	1
1:A:142:THR:CG2	1:A:196:LEU:HB3	0.42	2.44	19	1
1:A:112:PHE:CD2	1:A:229:TYR:HA	0.42	2.49	8	1
1:A:209:PHE:O	1:A:229:TYR:HB2	0.42	2.15	13	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:160:TYR:O	1:A:211:ILE:HA	0.42	2.15	2	1
1:A:119:GLU:O	1:A:123:GLN:HG3	0.42	2.15	16	2
1:A:111:LEU:HD13	1:A:111:LEU:H	0.41	1.73	10	1
1:A:183:CYS:HB3	1:A:188:ASP:OD2	0.41	2.16	9	1
1:A:125:LEU:CD1	1:A:131:MET:HG3	0.41	2.45	12	1
1:A:144:MET:H	1:A:194:ILE:HG22	0.41	1.76	2	2
1:A:111:LEU:HD12	1:A:111:LEU:N	0.41	2.31	11	3
1:A:136:GLU:CD	1:A:145:LYS:HD3	0.41	2.41	6	1
1:A:165:LEU:HD23	1:A:165:LEU:H	0.41	1.76	14	1
1:A:154:SER:CB	1:A:215:THR:HB	0.41	2.45	4	1
1:A:121:LEU:HD21	1:A:125:LEU:HD22	0.41	1.92	9	1
1:A:130:ALA:O	1:A:222:SER:HB2	0.41	2.15	9	1
1:A:106:TYR:CD1	1:A:137:TYR:HB2	0.41	2.51	8	1
1:A:128:GLN:O	1:A:130:ALA:N	0.41	2.54	17	1
1:A:129:LYS:O	1:A:150:VAL:HA	0.40	2.16	6	3
1:A:142:THR:HG22	1:A:196:LEU:CB	0.40	2.46	13	1
1:A:214:GLU:HG2	1:A:219:THR:HG23	0.40	1.92	16	1
1:A:131:MET:HE3	1:A:131:MET:C	0.40	2.41	20	1
1:A:118:LYS:O	1:A:122:MET:HG2	0.40	2.16	7	1
1:A:114:LEU:HD21	1:A:132:LEU:HD12	0.40	1.92	9	1
1:A:212:ARG:HD3	1:A:221:TRP:CZ2	0.40	2.51	12	1

6.3 Torsion angles ⓘ

6.3.1 Protein backbone ⓘ

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	131/139 (94%)	103±24 (78±18%)	17±4 (13±3%)	5±2 (4±1%)	4	32
All	All	2489/2780 (90%)	2052 (82%)	339 (14%)	98 (4%)	4	30

All 17 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	129	LYS	19

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Mol	Chain	Res	Type	Models (Total)
1	A	173	ASP	16
1	A	168	TRP	15
1	A	233	CYS	11
1	A	225	ASN	8
1	A	195	SER	6
1	A	106	TYR	6
1	A	198	PRO	5
1	A	204	GLY	2
1	A	115	PRO	2
1	A	224	ASN	2
1	A	207	VAL	1
1	A	228	ASN	1
1	A	201	GLN	1
1	A	229	TYR	1
1	A	206	LYS	1
1	A	139	PRO	1

6.3.2 Protein sidechains ⓘ

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	123/129 (95%)	106±27 (86±22%)	7±3 (5±2%)	21	71
All	All	2246/2580 (87%)	2111 (94%)	135 (6%)	19	68

All 26 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	128	GLN	18
1	A	172	TYR	18
1	A	111	LEU	11
1	A	108	LEU	8
1	A	118	LYS	8
1	A	131	MET	7
1	A	197	VAL	7
1	A	236	LYS	7
1	A	151	LEU	7

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Mol	Chain	Res	Type	Models (Total)
1	A	142	THR	7
1	A	231	LEU	5
1	A	165	LEU	4
1	A	202	LYS	4
1	A	201	GLN	3
1	A	206	LYS	3
1	A	196	LEU	3
1	A	149	ARG	2
1	A	158	LEU	2
1	A	160	TYR	2
1	A	137	TYR	2
1	A	175	LEU	2
1	A	132	LEU	1
1	A	123	GLN	1
1	A	189	GLN	1
1	A	106	TYR	1
1	A	234	GLN	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 ⓘ

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 82% for the well-defined parts and 81% 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	1524
Number of shifts mapped to atoms	1524
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	6

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	137	0.03 ± 0.14	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	131	-0.07 ± 0.14	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	128	0.03 ± 0.65	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 82%, i.e. 1476 atoms were assigned a chemical shift out of a possible 1803. 0 out of 23 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	518/654 (79%)	262/264 (99%)	132/264 (50%)	124/126 (98%)
Sidechain	899/985 (91%)	610/639 (95%)	275/314 (88%)	14/32 (44%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	59/164 (36%)	57/78 (73%)	0/82 (0%)	2/4 (50%)
Overall	1476/1803 (82%)	929/981 (95%)	407/660 (62%)	140/162 (86%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 81%, i.e. 1524 atoms were assigned a chemical shift out of a possible 1890. 0 out of 23 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	536/690 (78%)	271/279 (97%)	137/278 (49%)	128/133 (96%)
Sidechain	929/1028 (90%)	630/666 (95%)	285/329 (87%)	14/33 (42%)
Aromatic	59/172 (34%)	57/82 (70%)	0/84 (0%)	2/6 (33%)
Overall	1524/1890 (81%)	958/1027 (93%)	422/691 (61%)	144/172 (84%)

7.1.4 Statistically unusual chemical shifts ⓘ

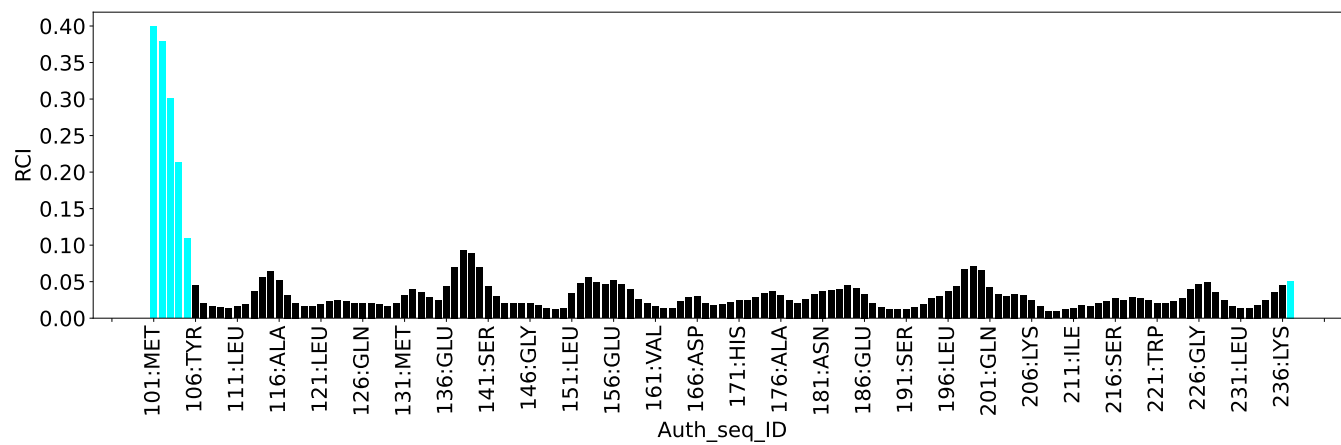
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	164	SER	HB3	1.45	2.49 – 5.20	-8.9
1	A	224	ASN	HB3	0.48	1.12 – 4.38	-7.0
1	A	212	ARG	HB3	0.05	0.43 – 3.11	-6.4
1	A	212	ARG	HG3	-0.16	0.15 – 2.94	-6.1
1	A	229	TYR	HB3	0.64	0.93 – 4.76	-5.8
1	A	164	SER	HB2	2.58	2.61 – 5.13	-5.1

7.1.5 Random Coil Index (RCI) plots ⓘ

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

Random coil index (RCI) for chain A:



8 NMR restraints analysis

8.1 Conformationally restricting restraints

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

Description	Value
Total distance restraints	2709
Intra-residue ($ i-j =0$)	596
Sequential ($ i-j =1$)	745
Medium range ($ i-j >1$ and $ i-j <5$)	313
Long range ($ i-j \geq 5$)	1055
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	19.5
Number of long range restraints per residue ¹	7.6

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)	20.2	0.2
0.2-0.5 (Medium)	0.7	0.32
>0.5 (Large)	None	None

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

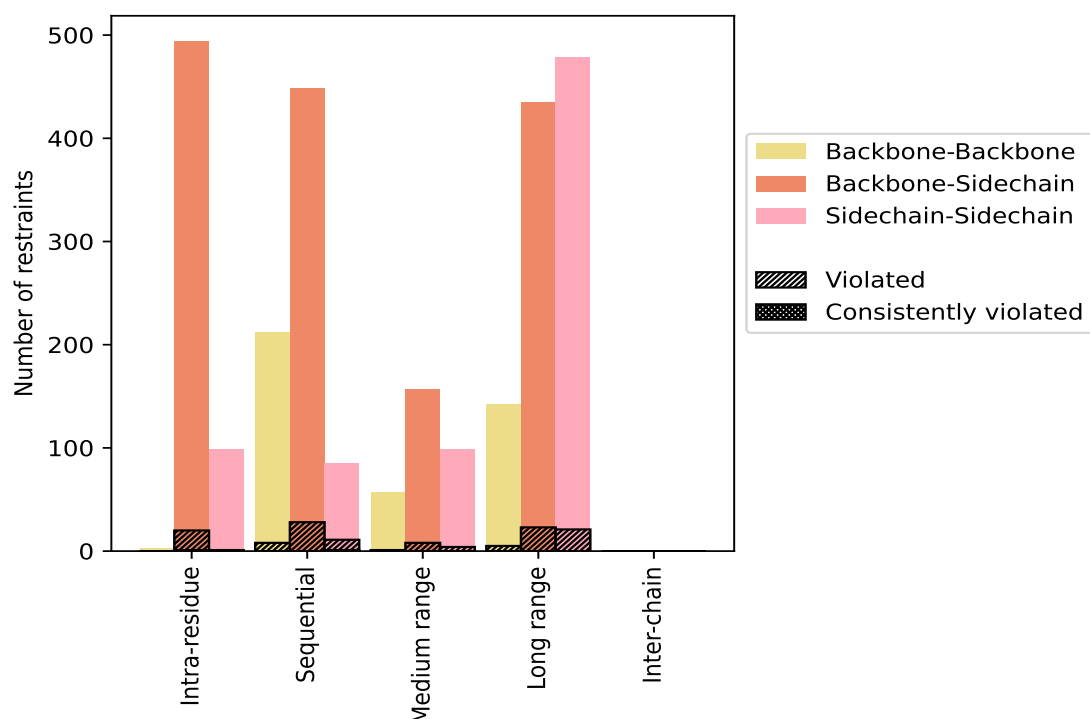
9.1 Summary of distance violations

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 ($i-j =0$)	596	22.0	21	3.5	0.8	1	0.2	0.0
Backbone-Backbone	3	0.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	494	18.2	20	4.0	0.7	1	0.2	0.0
Sidechain-Sidechain	99	3.7	1	1.0	0.0	0	0.0	0.0
Sequential ($i-j =1$)	745	27.5	47	6.3	1.7	1	0.1	0.0
Backbone-Backbone	212	7.8	8	3.8	0.3	0	0.0	0.0
Backbone-Sidechain	448	16.5	28	6.2	1.0	0	0.0	0.0
Sidechain-Sidechain	85	3.1	11	12.9	0.4	1	1.2	0.0
Medium range ($i-j >1$ & $i-j <5$)	313	11.6	13	4.2	0.5	0	0.0	0.0
Backbone-Backbone	57	2.1	1	1.8	0.0	0	0.0	0.0
Backbone-Sidechain	157	5.8	8	5.1	0.3	0	0.0	0.0
Sidechain-Sidechain	99	3.7	4	4.0	0.1	0	0.0	0.0
Long range ($i-j \geq 5$)	1055	38.9	49	4.6	1.8	0	0.0	0.0
Backbone-Backbone	142	5.2	5	3.5	0.2	0	0.0	0.0
Backbone-Sidechain	435	16.1	23	5.3	0.8	0	0.0	0.0
Sidechain-Sidechain	478	17.6	21	4.4	0.8	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	2709	100.0	130	4.8	4.8	2	0.1	0.1
Backbone-Backbone	414	15.3	14	3.4	0.5	0	0.0	0.0
Backbone-Sidechain	1534	56.6	79	5.1	2.9	1	0.1	0.0
Sidechain-Sidechain	761	28.1	37	4.9	1.4	1	0.1	0.0

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

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



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

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

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

Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	5	9	2	6	0	22	0.14	0.2	0.03	0.13
2	6	7	1	9	0	23	0.14	0.26	0.04	0.14
3	5	10	2	7	0	24	0.13	0.22	0.03	0.12
4	5	12	1	4	0	22	0.14	0.21	0.03	0.12
5	2	10	3	9	0	24	0.14	0.26	0.04	0.12
6	4	8	4	5	0	21	0.14	0.29	0.04	0.13
7	3	9	1	5	0	18	0.14	0.21	0.03	0.12
8	2	12	1	9	0	24	0.13	0.19	0.03	0.12
9	3	10	2	5	0	20	0.13	0.23	0.03	0.12
10	7	10	3	7	0	27	0.15	0.32	0.05	0.13

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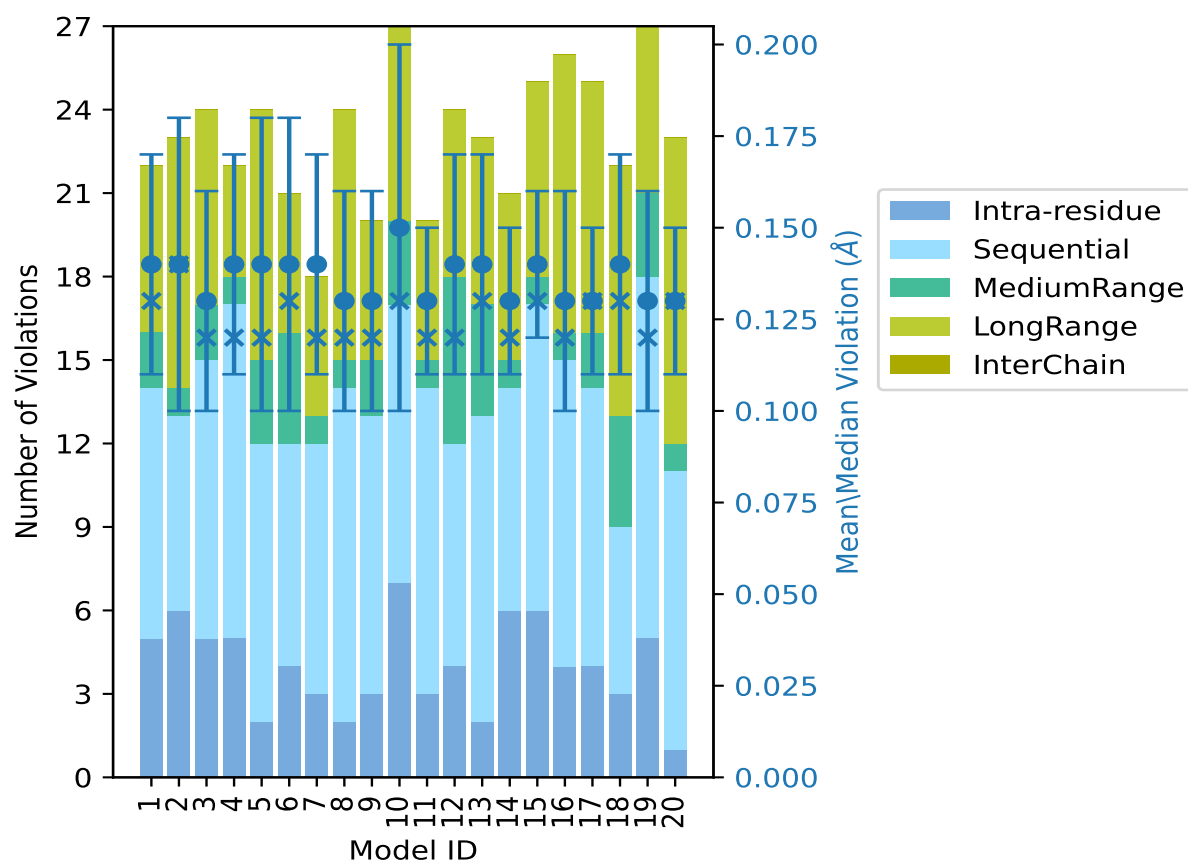
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
11	3	11	1	5	0	20	0.13	0.17	0.02	0.12
12	4	8	6	6	0	24	0.14	0.21	0.03	0.12
13	2	11	4	6	0	23	0.14	0.21	0.03	0.13
14	6	8	1	6	0	21	0.13	0.17	0.02	0.12
15	6	11	1	7	0	25	0.14	0.19	0.02	0.13
16	4	11	1	10	0	26	0.13	0.22	0.03	0.12
17	4	10	2	9	0	25	0.13	0.19	0.02	0.13
18	3	6	4	9	0	22	0.14	0.19	0.03	0.13
19	5	13	3	6	0	27	0.13	0.22	0.03	0.12
20	1	10	1	11	0	23	0.13	0.16	0.02	0.13

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

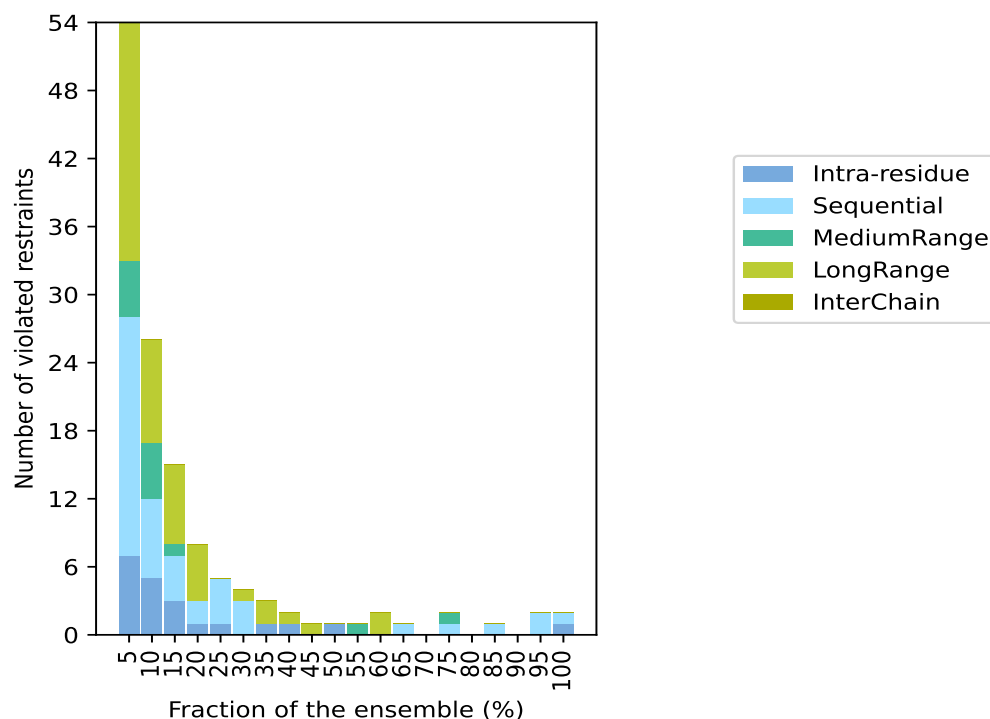
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 2579(IR:575, SQ:698, MR:300, LR:1006, IC:0) restraints are 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	%
7	21	5	21	0	54	1	5.0
5	7	5	9	0	26	2	10.0
3	4	1	7	0	15	3	15.0
1	2	0	5	0	8	4	20.0
1	4	0	0	0	5	5	25.0
0	3	0	1	0	4	6	30.0
1	0	0	2	0	3	7	35.0
1	0	0	1	0	2	8	40.0
0	0	0	1	0	1	9	45.0
1	0	0	0	0	1	10	50.0
0	0	1	0	0	1	11	55.0
0	0	0	2	0	2	12	60.0
0	1	0	0	0	1	13	65.0
0	0	0	0	0	0	14	70.0
0	1	1	0	0	2	15	75.0
0	0	0	0	0	0	16	80.0
0	1	0	0	0	1	17	85.0
0	0	0	0	0	0	18	90.0
0	2	0	0	0	2	19	95.0
1	1	0	0	0	2	20	100.0

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

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

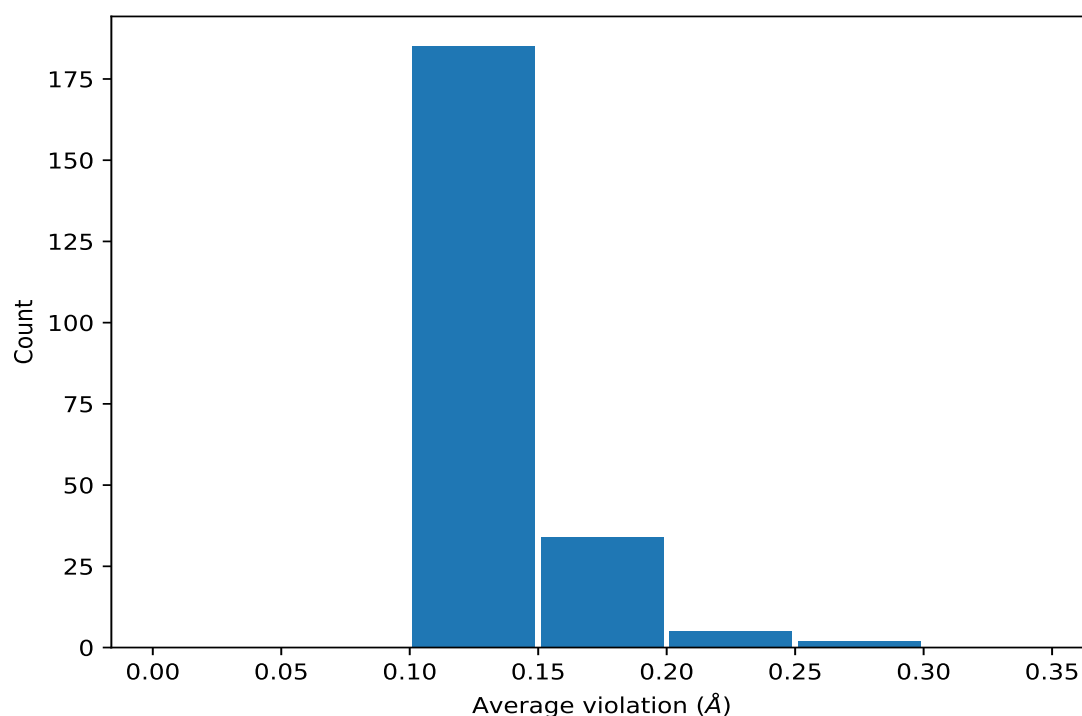
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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)
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	20	0.14	0.02	0.14
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	20	0.11	0.0	0.11
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	19	0.15	0.02	0.15
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	19	0.14	0.02	0.14
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	19	0.14	0.02	0.14
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	19	0.14	0.02	0.14
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	17	0.13	0.04	0.12
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	15	0.16	0.02	0.16
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	15	0.13	0.03	0.12
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	15	0.13	0.03	0.12
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	15	0.13	0.03	0.12
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	15	0.13	0.03	0.12
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	15	0.13	0.03	0.12
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	15	0.13	0.03	0.12
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	13	0.14	0.03	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	12	0.16	0.03	0.14

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	12	0.16	0.03	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	12	0.16	0.03	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	12	0.16	0.03	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	12	0.16	0.03	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	12	0.16	0.03	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	12	0.16	0.03	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	12	0.16	0.03	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	12	0.16	0.03	0.14
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	12	0.14	0.02	0.13
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	12	0.14	0.02	0.13
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	12	0.14	0.02	0.13
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	12	0.14	0.02	0.13
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	12	0.14	0.02	0.13
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	12	0.14	0.02	0.13
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	11	0.17	0.03	0.18
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	11	0.17	0.03	0.18
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	11	0.17	0.03	0.18
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	11	0.17	0.03	0.18
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	11	0.17	0.03	0.18
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	11	0.17	0.03	0.18
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	10	0.17	0.02	0.16
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	9	0.13	0.02	0.12
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	8	0.18	0.02	0.17
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	8	0.15	0.02	0.15
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	8	0.15	0.02	0.15
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD11	7	0.17	0.01	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD12	7	0.17	0.01	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD13	7	0.17	0.01	0.17
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD1	7	0.14	0.03	0.13
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD2	7	0.14	0.03	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG21	7	0.14	0.01	0.14
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG22	7	0.14	0.01	0.14
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG23	7	0.14	0.01	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG21	7	0.14	0.01	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG22	7	0.14	0.01	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG23	7	0.14	0.01	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG21	7	0.14	0.01	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG22	7	0.14	0.01	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG23	7	0.14	0.01	0.14
(1,315)	1:104:A:GLU:HA	1:105:A:GLU:H	6	0.14	0.02	0.15
(1,2699)	1:235:A:LYS:HB2	1:236:A:LYS:H	6	0.14	0.02	0.14
(1,2699)	1:235:A:LYS:HB3	1:236:A:LYS:H	6	0.14	0.02	0.14

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD1	6	0.12	0.01	0.12
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD2	6	0.12	0.01	0.12
(1,453)	1:218:A:GLY:H	1:219:A:THR:H	6	0.12	0.01	0.12
(1,823)	1:225:A:ASN:H	1:226:A:GLY:H	5	0.15	0.03	0.14
(1,1871)	1:236:A:LYS:HG2	1:237:A:GLU:H	5	0.13	0.02	0.12
(1,1871)	1:236:A:LYS:HG3	1:237:A:GLU:H	5	0.13	0.02	0.12
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG21	5	0.13	0.01	0.13
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG22	5	0.13	0.01	0.13
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG23	5	0.13	0.01	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG21	5	0.13	0.01	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG22	5	0.13	0.01	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG23	5	0.13	0.01	0.13
(1,2072)	1:111:A:LEU:HD11	1:112:A:PHE:H	5	0.13	0.02	0.12
(1,2072)	1:111:A:LEU:HD12	1:112:A:PHE:H	5	0.13	0.02	0.12
(1,2072)	1:111:A:LEU:HD13	1:112:A:PHE:H	5	0.13	0.02	0.12
(1,2072)	1:111:A:LEU:HD21	1:112:A:PHE:H	5	0.13	0.02	0.12
(1,2072)	1:111:A:LEU:HD22	1:112:A:PHE:H	5	0.13	0.02	0.12
(1,2072)	1:111:A:LEU:HD23	1:112:A:PHE:H	5	0.13	0.02	0.12
(1,1816)	1:108:A:LEU:H	1:108:A:LEU:HG	5	0.11	0.01	0.11
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE2	4	0.27	0.04	0.28
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE3	4	0.27	0.04	0.28
(1,2036)	1:105:A:GLU:HG2	1:106:A:TYR:HA	4	0.18	0.05	0.16
(1,2036)	1:105:A:GLU:HG3	1:106:A:TYR:HA	4	0.18	0.05	0.16
(1,2289)	1:149:A:ARG:HG2	1:188:A:ASP:H	4	0.16	0.03	0.17
(1,2289)	1:149:A:ARG:HG3	1:188:A:ASP:H	4	0.16	0.03	0.17
(1,1765)	1:160:A:TYR:HD1	1:212:A:ARG:HG3	4	0.13	0.02	0.12
(1,1765)	1:160:A:TYR:HD2	1:212:A:ARG:HG3	4	0.13	0.02	0.12
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD2	4	0.13	0.01	0.12
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD3	4	0.13	0.01	0.12
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG21	4	0.12	0.02	0.12
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG22	4	0.12	0.02	0.12
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG23	4	0.12	0.02	0.12
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG21	4	0.12	0.02	0.12
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG22	4	0.12	0.02	0.12
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG23	4	0.12	0.02	0.12
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD11	4	0.12	0.01	0.12
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD12	4	0.12	0.01	0.12
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD13	4	0.12	0.01	0.12
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD11	4	0.12	0.01	0.12
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD12	4	0.12	0.01	0.12
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD13	4	0.12	0.01	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG11	4	0.12	0.01	0.12

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG12	4	0.12	0.01	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG13	4	0.12	0.01	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG21	4	0.12	0.01	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG22	4	0.12	0.01	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG23	4	0.12	0.01	0.12
(1,2034)	1:105:A:GLU:HA	1:105:A:GLU:HG2	3	0.17	0.04	0.16
(1,2034)	1:105:A:GLU:HA	1:105:A:GLU:HG3	3	0.17	0.04	0.16
(1,2038)	1:106:A:TYR:H	1:106:A:TYR:HB2	3	0.16	0.02	0.17
(1,2038)	1:106:A:TYR:H	1:106:A:TYR:HB3	3	0.16	0.02	0.17
(1,1246)	1:205:A:SER:HB2	1:206:A:LYS:HE2	3	0.13	0.02	0.14
(1,1246)	1:205:A:SER:HB2	1:206:A:LYS:HE3	3	0.13	0.02	0.14
(1,1246)	1:205:A:SER:HB3	1:206:A:LYS:HE2	3	0.13	0.02	0.14
(1,1246)	1:205:A:SER:HB3	1:206:A:LYS:HE3	3	0.13	0.02	0.14
(1,1405)	1:137:A:TYR:HE1	1:234:A:GLN:HA	3	0.13	0.03	0.11
(1,1405)	1:137:A:TYR:HE2	1:234:A:GLN:HA	3	0.13	0.03	0.11
(1,1337)	1:143:A:SER:HA	1:194:A:ILE:HB	3	0.13	0.02	0.12
(1,383)	1:113:A:ASP:H	1:229:A:TYR:HA	3	0.12	0.01	0.12
(1,1055)	1:154:A:SER:HB3	1:157:A:LYS:HB3	3	0.12	0.01	0.11
(1,1577)	1:147:A:ILE:HB	1:191:A:SER:HB2	3	0.11	0.01	0.11
(1,2580)	1:202:A:LYS:HB2	1:202:A:LYS:HE2	3	0.11	0.0	0.11
(1,2580)	1:202:A:LYS:HB2	1:202:A:LYS:HE3	3	0.11	0.0	0.11
(1,2580)	1:202:A:LYS:HB3	1:202:A:LYS:HE2	3	0.11	0.0	0.11
(1,2580)	1:202:A:LYS:HB3	1:202:A:LYS:HE3	3	0.11	0.0	0.11
(1,371)	1:194:A:ILE:HA	1:195:A:SER:H	3	0.11	0.01	0.11
(1,526)	1:107:A:VAL:H	1:235:A:LYS:HA	3	0.11	0.01	0.11
(1,1062)	1:114:A:LEU:HD21	1:115:A:PRO:HD3	3	0.11	0.0	0.11
(1,1062)	1:114:A:LEU:HD22	1:115:A:PRO:HD3	3	0.11	0.0	0.11
(1,1062)	1:114:A:LEU:HD23	1:115:A:PRO:HD3	3	0.11	0.0	0.11
(1,968)	1:152:A:ASN:HD22	1:153:A:ILE:HA	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD11	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD12	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD13	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD11	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD12	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD13	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD11	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD12	3	0.11	0.0	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD13	3	0.11	0.0	0.11
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD11	3	0.1	0.0	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD12	3	0.1	0.0	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD13	3	0.1	0.0	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD21	3	0.1	0.0	0.1

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD22	3	0.1	0.0	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD23	3	0.1	0.0	0.1
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD21	2	0.2	0.01	0.2
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD22	2	0.2	0.01	0.2
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD23	2	0.2	0.01	0.2
(1,1009)	1:171:A:HIS:HA	1:172:A:TYR:HE1	2	0.2	0.02	0.2
(1,1009)	1:171:A:HIS:HA	1:172:A:TYR:HE2	2	0.2	0.02	0.2
(1,390)	1:105:A:GLU:HA	1:106:A:TYR:H	2	0.16	0.02	0.16
(1,269)	1:104:A:GLU:H	1:104:A:GLU:HB2	2	0.14	0.04	0.14
(1,269)	1:104:A:GLU:H	1:104:A:GLU:HB3	2	0.14	0.04	0.14
(1,2678)	1:231:A:LEU:HD11	1:234:A:GLN:H	2	0.13	0.01	0.13
(1,2678)	1:231:A:LEU:HD12	1:234:A:GLN:H	2	0.13	0.01	0.13
(1,2678)	1:231:A:LEU:HD13	1:234:A:GLN:H	2	0.13	0.01	0.13
(1,2678)	1:231:A:LEU:HD21	1:234:A:GLN:H	2	0.13	0.01	0.13
(1,2678)	1:231:A:LEU:HD22	1:234:A:GLN:H	2	0.13	0.01	0.13
(1,2678)	1:231:A:LEU:HD23	1:234:A:GLN:H	2	0.13	0.01	0.13
(1,2683)	1:232:A:VAL:HB	1:233:A:CYS:HB2	2	0.13	0.01	0.13
(1,2683)	1:232:A:VAL:HB	1:233:A:CYS:HB3	2	0.13	0.01	0.13
(1,842)	1:133:A:GLU:H	1:146:A:GLY:HA3	2	0.12	0.01	0.12
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG21	2	0.12	0.02	0.12
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG22	2	0.12	0.02	0.12
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG23	2	0.12	0.02	0.12
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG21	2	0.12	0.02	0.12
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG22	2	0.12	0.02	0.12
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG23	2	0.12	0.02	0.12
(1,1499)	1:196:A:LEU:HB2	1:200:A:TYR:HE1	2	0.12	0.02	0.12
(1,1499)	1:196:A:LEU:HB2	1:200:A:TYR:HE2	2	0.12	0.02	0.12
(1,1499)	1:196:A:LEU:HB3	1:200:A:TYR:HE1	2	0.12	0.02	0.12
(1,1499)	1:196:A:LEU:HB3	1:200:A:TYR:HE2	2	0.12	0.02	0.12
(1,2123)	1:121:A:LEU:HD11	1:125:A:LEU:HA	2	0.12	0.02	0.12
(1,2123)	1:121:A:LEU:HD12	1:125:A:LEU:HA	2	0.12	0.02	0.12
(1,2123)	1:121:A:LEU:HD13	1:125:A:LEU:HA	2	0.12	0.02	0.12
(1,2123)	1:121:A:LEU:HD21	1:125:A:LEU:HA	2	0.12	0.02	0.12
(1,2123)	1:121:A:LEU:HD22	1:125:A:LEU:HA	2	0.12	0.02	0.12
(1,2123)	1:121:A:LEU:HD23	1:125:A:LEU:HA	2	0.12	0.02	0.12
(1,1767)	1:120:A:ASP:HA	1:123:A:GLN:HB2	2	0.12	0.01	0.12
(1,1767)	1:120:A:ASP:HA	1:123:A:GLN:HB3	2	0.12	0.01	0.12
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG11	2	0.12	0.01	0.12
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG12	2	0.12	0.01	0.12
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG13	2	0.12	0.01	0.12
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG21	2	0.12	0.01	0.12
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG22	2	0.12	0.01	0.12

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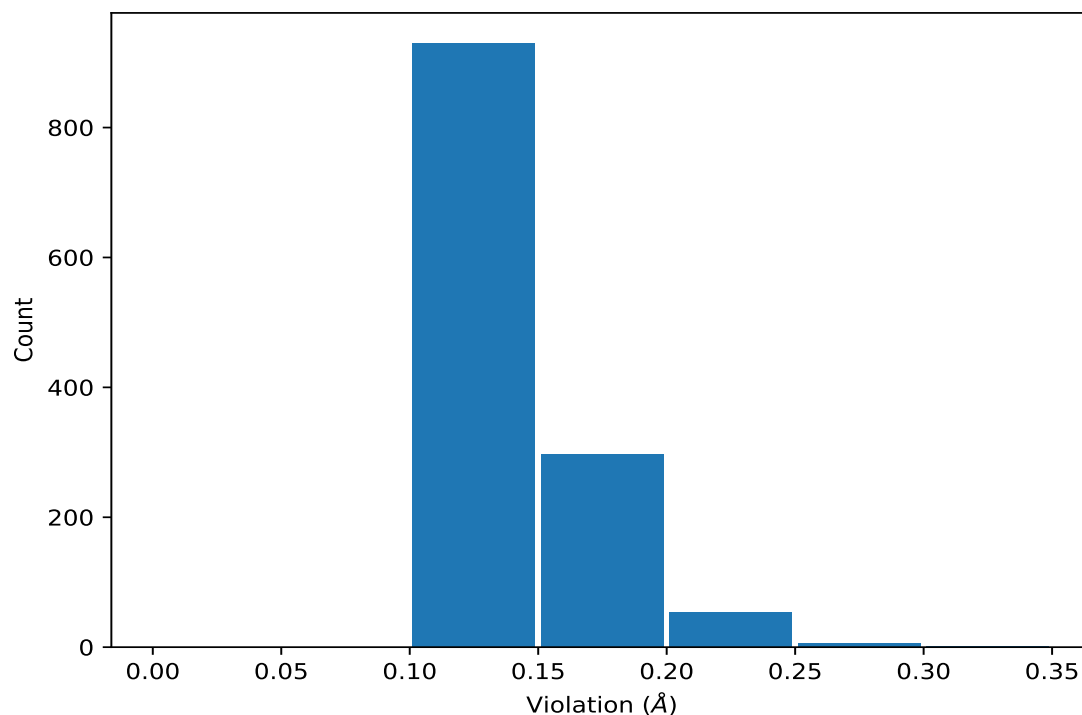
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG23	2	0.12	0.01	0.12
(1,1610)	1:156:A:GLU:H	1:156:A:GLU:HG3	2	0.12	0.0	0.12
(1,1841)	1:206:A:LYS:HD3	1:233:A:CYS:H	2	0.12	0.0	0.12
(1,2037)	1:105:A:GLU:HG2	1:236:A:LYS:HE2	2	0.12	0.0	0.12
(1,2037)	1:105:A:GLU:HG2	1:236:A:LYS:HE3	2	0.12	0.0	0.12
(1,2037)	1:105:A:GLU:HG3	1:236:A:LYS:HE2	2	0.12	0.0	0.12
(1,2037)	1:105:A:GLU:HG3	1:236:A:LYS:HE3	2	0.12	0.0	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG11	2	0.12	0.0	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG12	2	0.12	0.0	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG13	2	0.12	0.0	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG21	2	0.12	0.0	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG22	2	0.12	0.0	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG23	2	0.12	0.0	0.12
(1,2548)	1:193:A:LYS:HG2	1:194:A:ILE:HA	2	0.12	0.0	0.12
(1,2548)	1:193:A:LYS:HG3	1:194:A:ILE:HA	2	0.12	0.0	0.12
(1,95)	1:132:A:LEU:HD21	1:133:A:GLU:H	2	0.11	0.01	0.11
(1,95)	1:132:A:LEU:HD22	1:133:A:GLU:H	2	0.11	0.01	0.11
(1,95)	1:132:A:LEU:HD23	1:133:A:GLU:H	2	0.11	0.01	0.11
(1,113)	1:161:A:VAL:HG21	1:212:A:ARG:H	2	0.11	0.0	0.11
(1,113)	1:161:A:VAL:HG22	1:212:A:ARG:H	2	0.11	0.0	0.11
(1,113)	1:161:A:VAL:HG23	1:212:A:ARG:H	2	0.11	0.0	0.11
(1,1088)	1:236:A:LYS:HE2	1:237:A:GLU:HA	2	0.11	0.0	0.11
(1,1088)	1:236:A:LYS:HE3	1:237:A:GLU:HA	2	0.11	0.0	0.11
(1,1273)	1:172:A:TYR:HE1	1:174:A:ILE:HA	2	0.11	0.0	0.11
(1,1273)	1:172:A:TYR:HE2	1:174:A:ILE:HA	2	0.11	0.0	0.11
(1,2044)	1:106:A:TYR:HB2	1:236:A:LYS:H	2	0.11	0.01	0.11
(1,2044)	1:106:A:TYR:HB3	1:236:A:LYS:H	2	0.11	0.01	0.11
(1,238)	1:173:A:ASP:H	1:173:A:ASP:HB3	2	0.11	0.0	0.11
(1,581)	1:128:A:GLN:H	1:128:A:GLN:HG3	2	0.11	0.0	0.11
(1,1747)	1:182:A:SER:H	1:183:A:CYS:HB2	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD11	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD12	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD13	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD21	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD22	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD23	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD11	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD12	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD13	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD21	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD22	2	0.11	0.0	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD23	2	0.11	0.0	0.11

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)
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE2	10	0.32
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE3	10	0.32
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE2	6	0.29
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE3	6	0.29
(1,2036)	1:105:A:GLU:HG2	1:106:A:TYR:HA	5	0.26
(1,2036)	1:105:A:GLU:HG3	1:106:A:TYR:HA	5	0.26
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE2	2	0.26

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE3	2	0.26
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	9	0.23
(1,2034)	1:105:A:GLU:HA	1:105:A:GLU:HG2	16	0.22
(1,2034)	1:105:A:GLU:HA	1:105:A:GLU:HG3	16	0.22
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	19	0.22
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	19	0.22
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	19	0.22
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	19	0.22
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	19	0.22
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	19	0.22
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	19	0.22
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	19	0.22
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	19	0.22
(1,1099)	1:156:A:GLU:HA	1:156:A:GLU:HG3	3	0.22
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD21	12	0.21
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD22	12	0.21
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD23	12	0.21
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	13	0.21
(1,1009)	1:171:A:HIS:HA	1:172:A:TYR:HE1	5	0.21
(1,1009)	1:171:A:HIS:HA	1:172:A:TYR:HE2	5	0.21
(1,823)	1:225:A:ASN:H	1:226:A:GLY:H	7	0.21
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	13	0.21
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	4	0.21
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	5	0.2
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	5	0.2
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	5	0.2
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	5	0.2
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	5	0.2
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	5	0.2
(1,2289)	1:149:A:ARG:HG2	1:188:A:ASP:H	10	0.2
(1,2289)	1:149:A:ARG:HG3	1:188:A:ASP:H	10	0.2
(1,2036)	1:105:A:GLU:HG2	1:106:A:TYR:HA	4	0.2
(1,2036)	1:105:A:GLU:HG3	1:106:A:TYR:HA	4	0.2
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	1	0.2
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	1	0.2
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	1	0.2
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	1	0.2
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	1	0.2
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	1	0.2
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	1	0.2
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	1	0.2
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	1	0.2

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	4	0.2
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	4	0.2
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	4	0.2
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	4	0.2
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	4	0.2
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	4	0.2
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	4	0.2
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	4	0.2
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	4	0.2
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE2	19	0.2
(1,1547)	1:145:A:LYS:HA	1:145:A:LYS:HE3	19	0.2
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	5	0.2
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	7	0.2
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	2	0.19
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	2	0.19
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	2	0.19
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	2	0.19
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	2	0.19
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	2	0.19
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	8	0.19
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	8	0.19
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	8	0.19
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	8	0.19
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	8	0.19
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	8	0.19
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	18	0.19
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	18	0.19
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	18	0.19
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	18	0.19
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	18	0.19
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	18	0.19
(1,2289)	1:149:A:ARG:HG2	1:188:A:ASP:H	8	0.19
(1,2289)	1:149:A:ARG:HG3	1:188:A:ASP:H	8	0.19
(1,2038)	1:106:A:TYR:H	1:106:A:TYR:HB2	15	0.19
(1,2038)	1:106:A:TYR:H	1:106:A:TYR:HB3	15	0.19
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD21	4	0.19
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD22	4	0.19
(1,1873)	1:158:A:LEU:HA	1:158:A:LEU:HD23	4	0.19
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	3	0.19
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	10	0.19
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	10	0.19
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	12	0.19

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,390)	1:105:A:GLU:HA	1:106:A:TYR:H	12	0.19
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	2	0.19
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	6	0.19
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	17	0.19
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	10	0.18
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	10	0.18
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	10	0.18
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	10	0.18
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	10	0.18
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	10	0.18
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	12	0.18
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	12	0.18
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	12	0.18
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	12	0.18
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	12	0.18
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	12	0.18
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	13	0.18
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	13	0.18
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	13	0.18
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	13	0.18
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	13	0.18
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	13	0.18
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	18	0.18
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	18	0.18
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	1	0.18
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	1	0.18
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	1	0.18
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	1	0.18
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	1	0.18
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	1	0.18
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	2	0.18
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	2	0.18
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	2	0.18
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	2	0.18
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	2	0.18
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	2	0.18
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD1	9	0.18
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD2	9	0.18
(1,1405)	1:137:A:TYR:HE1	1:234:A:GLN:HA	12	0.18
(1,1405)	1:137:A:TYR:HE2	1:234:A:GLN:HA	12	0.18
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD11	1	0.18
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD12	1	0.18

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD13	1	0.18
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD11	15	0.18
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD12	15	0.18
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD13	15	0.18
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	18	0.18
(1,1009)	1:171:A:HIS:HA	1:172:A:TYR:HE1	12	0.18
(1,1009)	1:171:A:HIS:HA	1:172:A:TYR:HE2	12	0.18
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	10	0.18
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	10	0.18
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	1	0.18
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	3	0.18
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	11	0.17
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	11	0.17
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	11	0.17
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	11	0.17
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	11	0.17
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	11	0.17
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	3	0.17
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	3	0.17
(1,2045)	1:107:A:VAL:H	1:235:A:LYS:HG2	18	0.17
(1,2045)	1:107:A:VAL:H	1:235:A:LYS:HG3	18	0.17
(1,2038)	1:106:A:TYR:H	1:106:A:TYR:HB2	14	0.17
(1,2038)	1:106:A:TYR:H	1:106:A:TYR:HB3	14	0.17
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	18	0.17
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	18	0.17
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	18	0.17
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	18	0.17
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	18	0.17
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	18	0.17
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	17	0.17
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	17	0.17
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	17	0.17
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	17	0.17
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	17	0.17
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	17	0.17
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	17	0.17
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	17	0.17
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	17	0.17
(1,1765)	1:160:A:TYR:HD1	1:212:A:ARG:HG3	2	0.17
(1,1765)	1:160:A:TYR:HD2	1:212:A:ARG:HG3	2	0.17
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD1	18	0.17
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD2	18	0.17

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD11	6	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD12	6	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD13	6	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD11	18	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD12	18	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD13	18	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD11	19	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD12	19	0.17
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD13	19	0.17
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	6	0.17
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	9	0.17
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	12	0.17
(1,1214)	1:143:A:SER:HB2	1:196:A:LEU:HA	15	0.17
(1,1214)	1:143:A:SER:HB3	1:196:A:LEU:HA	15	0.17
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	11	0.17
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	12	0.17
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	17	0.17
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	19	0.17
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	15	0.17
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	19	0.17
(1,269)	1:104:A:GLU:H	1:104:A:GLU:HB2	9	0.17
(1,269)	1:104:A:GLU:H	1:104:A:GLU:HB3	9	0.17
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	16	0.17
(1,2699)	1:235:A:LYS:HB2	1:236:A:LYS:H	20	0.16
(1,2699)	1:235:A:LYS:HB3	1:236:A:LYS:H	20	0.16
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	11	0.16
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	11	0.16
(1,2034)	1:105:A:GLU:HA	1:105:A:GLU:HG2	10	0.16
(1,2034)	1:105:A:GLU:HA	1:105:A:GLU:HG3	10	0.16
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	6	0.16
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	6	0.16
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	6	0.16
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	6	0.16
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	6	0.16
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	6	0.16
(1,1871)	1:236:A:LYS:HG2	1:237:A:GLU:H	20	0.16
(1,1871)	1:236:A:LYS:HG3	1:237:A:GLU:H	20	0.16
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD11	7	0.16
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD12	7	0.16
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD13	7	0.16
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD11	14	0.16
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD12	14	0.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1388)	1:158:A:LEU:HA	1:158:A:LEU:HD13	14	0.16
(1,1337)	1:143:A:SER:HA	1:194:A:ILE:HB	20	0.16
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	8	0.16
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	15	0.16
(1,1259)	1:103:A:THR:HA	1:104:A:GLU:H	8	0.16
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	4	0.16
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	4	0.16
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	4	0.16
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	5	0.16
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	5	0.16
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	5	0.16
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	4	0.16
(1,1100)	1:236:A:LYS:HE2	1:237:A:GLU:HG2	2	0.16
(1,1100)	1:236:A:LYS:HE2	1:237:A:GLU:HG3	2	0.16
(1,1100)	1:236:A:LYS:HE3	1:237:A:GLU:HG2	2	0.16
(1,1100)	1:236:A:LYS:HE3	1:237:A:GLU:HG3	2	0.16
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	2	0.16
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	4	0.16
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	16	0.16
(1,823)	1:225:A:ASN:H	1:226:A:GLY:H	15	0.16
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	3	0.16
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	4	0.16
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	16	0.16
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	20	0.16
(1,315)	1:104:A:GLU:HA	1:105:A:GLU:H	19	0.16
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	10	0.16
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	13	0.16
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	14	0.16
(1,190)	1:160:A:TYR:HD1	1:174:A:ILE:H	8	0.16
(1,190)	1:160:A:TYR:HD2	1:174:A:ILE:H	8	0.16
(1,2699)	1:235:A:LYS:HB2	1:236:A:LYS:H	13	0.15
(1,2699)	1:235:A:LYS:HB3	1:236:A:LYS:H	13	0.15
(1,2699)	1:235:A:LYS:HB2	1:236:A:LYS:H	19	0.15
(1,2699)	1:235:A:LYS:HB3	1:236:A:LYS:H	19	0.15
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	17	0.15
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	17	0.15
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	17	0.15
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	17	0.15
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	17	0.15
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	17	0.15
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	2	0.15
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	2	0.15

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	17	0.15
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	17	0.15
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	14	0.15
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	14	0.15
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	14	0.15
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	14	0.15
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	14	0.15
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	14	0.15
(1,2289)	1:149:A:ARG:HG2	1:188:A:ASP:H	15	0.15
(1,2289)	1:149:A:ARG:HG3	1:188:A:ASP:H	15	0.15
(1,2217)	1:134:A:SER:HB2	1:135:A:THR:H	14	0.15
(1,2217)	1:134:A:SER:HB3	1:135:A:THR:H	14	0.15
(1,2072)	1:111:A:LEU:HD11	1:112:A:PHE:H	11	0.15
(1,2072)	1:111:A:LEU:HD12	1:112:A:PHE:H	11	0.15
(1,2072)	1:111:A:LEU:HD13	1:112:A:PHE:H	11	0.15
(1,2072)	1:111:A:LEU:HD21	1:112:A:PHE:H	11	0.15
(1,2072)	1:111:A:LEU:HD22	1:112:A:PHE:H	11	0.15
(1,2072)	1:111:A:LEU:HD23	1:112:A:PHE:H	11	0.15
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	9	0.15
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	9	0.15
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	9	0.15
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	9	0.15
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	9	0.15
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	9	0.15
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	20	0.15
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	20	0.15
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	20	0.15
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	20	0.15
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	20	0.15
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	20	0.15
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG21	14	0.15
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG22	14	0.15
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG23	14	0.15
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG21	14	0.15
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG22	14	0.15
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG23	14	0.15
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG21	14	0.15
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG22	14	0.15
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG23	14	0.15
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG21	16	0.15
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG22	16	0.15
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG23	16	0.15

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG21	16	0.15
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG22	16	0.15
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG23	16	0.15
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG21	16	0.15
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG22	16	0.15
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG23	16	0.15
(1,1871)	1:236:A:LYS:HG2	1:237:A:GLU:H	15	0.15
(1,1871)	1:236:A:LYS:HG3	1:237:A:GLU:H	15	0.15
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	8	0.15
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	8	0.15
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	8	0.15
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	8	0.15
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	8	0.15
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	8	0.15
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	8	0.15
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	8	0.15
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	8	0.15
(1,1499)	1:196:A:LEU:HB2	1:200:A:TYR:HE1	6	0.15
(1,1499)	1:196:A:LEU:HB2	1:200:A:TYR:HE2	6	0.15
(1,1499)	1:196:A:LEU:HB3	1:200:A:TYR:HE1	6	0.15
(1,1499)	1:196:A:LEU:HB3	1:200:A:TYR:HE2	6	0.15
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	12	0.15
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	17	0.15
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	19	0.15
(1,1280)	1:151:A:LEU:HG	1:187:A:THR:HA	10	0.15
(1,1246)	1:205:A:SER:HB2	1:206:A:LYS:HE2	11	0.15
(1,1246)	1:205:A:SER:HB2	1:206:A:LYS:HE3	11	0.15
(1,1246)	1:205:A:SER:HB3	1:206:A:LYS:HE2	11	0.15
(1,1246)	1:205:A:SER:HB3	1:206:A:LYS:HE3	11	0.15
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	1	0.15
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	1	0.15
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	1	0.15
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	3	0.15
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	3	0.15
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	3	0.15
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	17	0.15
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	17	0.15
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	17	0.15
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	20	0.15
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	20	0.15
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	20	0.15
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	11	0.15

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	3	0.15
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	7	0.15
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	14	0.15
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	15	0.15
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	20	0.15
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	2	0.15
(1,546)	1:149:A:ARG:HD3	1:150:A:VAL:H	10	0.15
(1,541)	1:149:A:ARG:HD2	1:150:A:VAL:H	4	0.15
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	1	0.15
(1,513)	1:109:A:SER:H	1:109:A:SER:HB3	2	0.15
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	2	0.15
(1,355)	1:207:A:VAL:H	1:207:A:VAL:HB	1	0.15
(1,315)	1:104:A:GLU:HA	1:105:A:GLU:H	4	0.15
(1,315)	1:104:A:GLU:HA	1:105:A:GLU:H	11	0.15
(1,315)	1:104:A:GLU:HA	1:105:A:GLU:H	15	0.15
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	19	0.15
(1,2683)	1:232:A:VAL:HB	1:233:A:CYS:HB2	18	0.14
(1,2683)	1:232:A:VAL:HB	1:233:A:CYS:HB3	18	0.14
(1,2678)	1:231:A:LEU:HD11	1:234:A:GLN:H	5	0.14
(1,2678)	1:231:A:LEU:HD12	1:234:A:GLN:H	5	0.14
(1,2678)	1:231:A:LEU:HD13	1:234:A:GLN:H	5	0.14
(1,2678)	1:231:A:LEU:HD21	1:234:A:GLN:H	5	0.14
(1,2678)	1:231:A:LEU:HD22	1:234:A:GLN:H	5	0.14
(1,2678)	1:231:A:LEU:HD23	1:234:A:GLN:H	5	0.14
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	3	0.14
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	3	0.14
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	3	0.14
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	3	0.14
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	3	0.14
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	3	0.14
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG21	13	0.14
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG22	13	0.14
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG23	13	0.14
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG21	13	0.14
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG22	13	0.14
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG23	13	0.14
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	7	0.14
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	7	0.14
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	7	0.14
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	7	0.14
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	7	0.14
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	7	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	17	0.14
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	17	0.14
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	17	0.14
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	17	0.14
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	17	0.14
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	17	0.14
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD11	20	0.14
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD12	20	0.14
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD13	20	0.14
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD11	20	0.14
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD12	20	0.14
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD13	20	0.14
(1,2123)	1:121:A:LEU:HD11	1:125:A:LEU:HA	12	0.14
(1,2123)	1:121:A:LEU:HD12	1:125:A:LEU:HA	12	0.14
(1,2123)	1:121:A:LEU:HD13	1:125:A:LEU:HA	12	0.14
(1,2123)	1:121:A:LEU:HD21	1:125:A:LEU:HA	12	0.14
(1,2123)	1:121:A:LEU:HD22	1:125:A:LEU:HA	12	0.14
(1,2123)	1:121:A:LEU:HD23	1:125:A:LEU:HA	12	0.14
(1,2072)	1:111:A:LEU:HD11	1:112:A:PHE:H	8	0.14
(1,2072)	1:111:A:LEU:HD12	1:112:A:PHE:H	8	0.14
(1,2072)	1:111:A:LEU:HD13	1:112:A:PHE:H	8	0.14
(1,2072)	1:111:A:LEU:HD21	1:112:A:PHE:H	8	0.14
(1,2072)	1:111:A:LEU:HD22	1:112:A:PHE:H	8	0.14
(1,2072)	1:111:A:LEU:HD23	1:112:A:PHE:H	8	0.14
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG21	2	0.14
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG22	2	0.14
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG23	2	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG21	2	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG22	2	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG23	2	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG21	2	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG22	2	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG23	2	0.14
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG21	20	0.14
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG22	20	0.14
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG23	20	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG21	20	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG22	20	0.14
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG23	20	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG21	20	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG22	20	0.14
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG23	20	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	11	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	11	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	11	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	11	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	11	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	11	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	11	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	11	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	11	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	13	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	13	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	13	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	13	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	13	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	13	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	13	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	13	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	13	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	16	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	16	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	16	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	16	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	16	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	16	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	16	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	16	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	16	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	20	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	20	0.14
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	20	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	20	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	20	0.14
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	20	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	20	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	20	0.14
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	20	0.14
(1,1590)	1:208:A:GLU:HG2	1:230:A:THR:HG21	20	0.14
(1,1590)	1:208:A:GLU:HG2	1:230:A:THR:HG22	20	0.14
(1,1590)	1:208:A:GLU:HG2	1:230:A:THR:HG23	20	0.14
(1,1246)	1:205:A:SER:HB2	1:206:A:LYS:HE2	10	0.14
(1,1246)	1:205:A:SER:HB2	1:206:A:LYS:HE3	10	0.14
(1,1246)	1:205:A:SER:HB3	1:206:A:LYS:HE2	10	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1246)	1:205:A:SER:HB3	1:206:A:LYS:HE3	10	0.14
(1,1232)	1:154:A:SER:HB2	1:157:A:LYS:HG2	12	0.14
(1,1232)	1:154:A:SER:HB2	1:157:A:LYS:HG3	12	0.14
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	6	0.14
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	6	0.14
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	6	0.14
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	7	0.14
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	7	0.14
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	7	0.14
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	8	0.14
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	8	0.14
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	8	0.14
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	11	0.14
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	11	0.14
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	11	0.14
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	13	0.14
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	13	0.14
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	13	0.14
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	7	0.14
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	13	0.14
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	16	0.14
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG21	6	0.14
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG22	6	0.14
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG23	6	0.14
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG21	6	0.14
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG22	6	0.14
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG23	6	0.14
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG21	7	0.14
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG22	7	0.14
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG23	7	0.14
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG21	7	0.14
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG22	7	0.14
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG23	7	0.14
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD2	8	0.14
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD3	8	0.14
(1,1092)	1:160:A:TYR:HD1	1:173:A:ASP:HB2	17	0.14
(1,1092)	1:160:A:TYR:HD2	1:173:A:ASP:HB2	17	0.14
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	1	0.14
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	9	0.14
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	13	0.14
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	1	0.14
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	3	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	5	0.14
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	6	0.14
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	9	0.14
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	13	0.14
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	14	0.14
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	15	0.14
(1,823)	1:225:A:ASN:H	1:226:A:GLY:H	5	0.14
(1,390)	1:105:A:GLU:HA	1:106:A:TYR:H	3	0.14
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	14	0.14
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	8	0.14
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	9	0.14
(1,166)	1:206:A:LYS:HD2	1:233:A:CYS:H	16	0.14
(1,2699)	1:235:A:LYS:HB2	1:236:A:LYS:H	3	0.13
(1,2699)	1:235:A:LYS:HB3	1:236:A:LYS:H	3	0.13
(1,2677)	1:231:A:LEU:HD11	1:233:A:CYS:H	13	0.13
(1,2677)	1:231:A:LEU:HD12	1:233:A:CYS:H	13	0.13
(1,2677)	1:231:A:LEU:HD13	1:233:A:CYS:H	13	0.13
(1,2677)	1:231:A:LEU:HD21	1:233:A:CYS:H	13	0.13
(1,2677)	1:231:A:LEU:HD22	1:233:A:CYS:H	13	0.13
(1,2677)	1:231:A:LEU:HD23	1:233:A:CYS:H	13	0.13
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG11	17	0.13
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG12	17	0.13
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG13	17	0.13
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG21	17	0.13
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG22	17	0.13
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG23	17	0.13
(1,2563)	1:199:A:PRO:HA	1:235:A:LYS:HB2	16	0.13
(1,2563)	1:199:A:PRO:HA	1:235:A:LYS:HB3	16	0.13
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	18	0.13
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	18	0.13
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	18	0.13
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	18	0.13
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	18	0.13
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	18	0.13
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG21	1	0.13
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG22	1	0.13
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG23	1	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG21	1	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG22	1	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG23	1	0.13
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG21	20	0.13
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG22	20	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG23	20	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG21	20	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG22	20	0.13
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG23	20	0.13
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	9	0.13
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	9	0.13
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG11	1	0.13
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG12	1	0.13
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG13	1	0.13
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG21	1	0.13
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG22	1	0.13
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG23	1	0.13
(1,2054)	1:109:A:SER:H	1:232:A:VAL:HG11	18	0.13
(1,2054)	1:109:A:SER:H	1:232:A:VAL:HG12	18	0.13
(1,2054)	1:109:A:SER:H	1:232:A:VAL:HG13	18	0.13
(1,2054)	1:109:A:SER:H	1:232:A:VAL:HG21	18	0.13
(1,2054)	1:109:A:SER:H	1:232:A:VAL:HG22	18	0.13
(1,2054)	1:109:A:SER:H	1:232:A:VAL:HG23	18	0.13
(1,2038)	1:106:A:TYR:H	1:106:A:TYR:HB2	10	0.13
(1,2038)	1:106:A:TYR:H	1:106:A:TYR:HB3	10	0.13
(1,2036)	1:105:A:GLU:HG2	1:106:A:TYR:HA	9	0.13
(1,2036)	1:105:A:GLU:HG3	1:106:A:TYR:HA	9	0.13
(1,2036)	1:105:A:GLU:HG2	1:106:A:TYR:HA	17	0.13
(1,2036)	1:105:A:GLU:HG3	1:106:A:TYR:HA	17	0.13
(1,2034)	1:105:A:GLU:HA	1:105:A:GLU:HG2	2	0.13
(1,2034)	1:105:A:GLU:HA	1:105:A:GLU:HG3	2	0.13
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	1	0.13
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	1	0.13
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	1	0.13
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	1	0.13
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	1	0.13
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	1	0.13
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	15	0.13
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	15	0.13
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	15	0.13
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	15	0.13
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	15	0.13
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	15	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG21	8	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG22	8	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG23	8	0.13
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG21	8	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG22	8	0.13
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG23	8	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG21	8	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG22	8	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG23	8	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG21	13	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG22	13	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG23	13	0.13
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG21	13	0.13
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG22	13	0.13
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG23	13	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG21	13	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG22	13	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG23	13	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG21	18	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG22	18	0.13
(1,1887)	1:151:A:LEU:HD11	1:187:A:THR:HG23	18	0.13
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG21	18	0.13
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG22	18	0.13
(1,1887)	1:151:A:LEU:HD12	1:187:A:THR:HG23	18	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG21	18	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG22	18	0.13
(1,1887)	1:151:A:LEU:HD13	1:187:A:THR:HG23	18	0.13
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	14	0.13
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	14	0.13
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	14	0.13
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	14	0.13
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	14	0.13
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	14	0.13
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	14	0.13
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	14	0.13
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	14	0.13
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD1	10	0.13
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD2	10	0.13
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD1	19	0.13
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD2	19	0.13
(1,1767)	1:120:A:ASP:HA	1:123:A:GLN:HB2	13	0.13
(1,1767)	1:120:A:ASP:HA	1:123:A:GLN:HB3	13	0.13
(1,1765)	1:160:A:TYR:HD1	1:212:A:ARG:HG3	3	0.13
(1,1765)	1:160:A:TYR:HD2	1:212:A:ARG:HG3	3	0.13
(1,1577)	1:147:A:ILE:HB	1:191:A:SER:HB2	18	0.13
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD1	12	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD2	12	0.13
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD1	15	0.13
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD2	15	0.13
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD1	16	0.13
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD2	16	0.13
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD1	19	0.13
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD2	19	0.13
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	2	0.13
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	2	0.13
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	2	0.13
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	15	0.13
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	15	0.13
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	15	0.13
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	16	0.13
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	16	0.13
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	16	0.13
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	18	0.13
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	18	0.13
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	18	0.13
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	6	0.13
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG21	5	0.13
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG22	5	0.13
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG23	5	0.13
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG21	5	0.13
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG22	5	0.13
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG23	5	0.13
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD2	7	0.13
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD3	7	0.13
(1,1055)	1:154:A:SER:HB3	1:157:A:LYS:HB3	19	0.13
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	5	0.13
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	18	0.13
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	4	0.13
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	8	0.13
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	18	0.13
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	19	0.13
(1,842)	1:133:A:GLU:H	1:146:A:GLY:HA3	6	0.13
(1,630)	1:142:A:THR:HG21	1:143:A:SER:H	16	0.13
(1,630)	1:142:A:THR:HG22	1:143:A:SER:H	16	0.13
(1,630)	1:142:A:THR:HG23	1:143:A:SER:H	16	0.13
(1,453)	1:218:A:GLY:H	1:219:A:THR:H	16	0.13
(1,398)	1:130:A:ALA:H	1:150:A:VAL:HA	17	0.13
(1,383)	1:113:A:ASP:H	1:229:A:TYR:HA	17	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,340)	1:160:A:TYR:HD1	1:161:A:VAL:H	17	0.13
(1,340)	1:160:A:TYR:HD2	1:161:A:VAL:H	17	0.13
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	20	0.13
(1,2699)	1:235:A:LYS:HB2	1:236:A:LYS:H	6	0.12
(1,2699)	1:235:A:LYS:HB3	1:236:A:LYS:H	6	0.12
(1,2699)	1:235:A:LYS:HB2	1:236:A:LYS:H	11	0.12
(1,2699)	1:235:A:LYS:HB3	1:236:A:LYS:H	11	0.12
(1,2683)	1:232:A:VAL:HB	1:233:A:CYS:HB2	14	0.12
(1,2683)	1:232:A:VAL:HB	1:233:A:CYS:HB3	14	0.12
(1,2678)	1:231:A:LEU:HD11	1:234:A:GLN:H	18	0.12
(1,2678)	1:231:A:LEU:HD12	1:234:A:GLN:H	18	0.12
(1,2678)	1:231:A:LEU:HD13	1:234:A:GLN:H	18	0.12
(1,2678)	1:231:A:LEU:HD21	1:234:A:GLN:H	18	0.12
(1,2678)	1:231:A:LEU:HD22	1:234:A:GLN:H	18	0.12
(1,2678)	1:231:A:LEU:HD23	1:234:A:GLN:H	18	0.12
(1,2666)	1:230:A:THR:HB	1:231:A:LEU:HD11	14	0.12
(1,2666)	1:230:A:THR:HB	1:231:A:LEU:HD12	14	0.12
(1,2666)	1:230:A:THR:HB	1:231:A:LEU:HD13	14	0.12
(1,2666)	1:230:A:THR:HB	1:231:A:LEU:HD21	14	0.12
(1,2666)	1:230:A:THR:HB	1:231:A:LEU:HD22	14	0.12
(1,2666)	1:230:A:THR:HB	1:231:A:LEU:HD23	14	0.12
(1,2580)	1:202:A:LYS:HB2	1:202:A:LYS:HE2	17	0.12
(1,2580)	1:202:A:LYS:HB2	1:202:A:LYS:HE3	17	0.12
(1,2580)	1:202:A:LYS:HB3	1:202:A:LYS:HE2	17	0.12
(1,2580)	1:202:A:LYS:HB3	1:202:A:LYS:HE3	17	0.12
(1,2551)	1:195:A:SER:HA	1:196:A:LEU:HD11	19	0.12
(1,2551)	1:195:A:SER:HA	1:196:A:LEU:HD12	19	0.12
(1,2551)	1:195:A:SER:HA	1:196:A:LEU:HD13	19	0.12
(1,2551)	1:195:A:SER:HA	1:196:A:LEU:HD21	19	0.12
(1,2551)	1:195:A:SER:HA	1:196:A:LEU:HD22	19	0.12
(1,2551)	1:195:A:SER:HA	1:196:A:LEU:HD23	19	0.12
(1,2548)	1:193:A:LYS:HG2	1:194:A:ILE:HA	20	0.12
(1,2548)	1:193:A:LYS:HG3	1:194:A:ILE:HA	20	0.12
(1,2491)	1:179:A:VAL:HG11	1:182:A:SER:HA	6	0.12
(1,2491)	1:179:A:VAL:HG12	1:182:A:SER:HA	6	0.12
(1,2491)	1:179:A:VAL:HG13	1:182:A:SER:HA	6	0.12
(1,2491)	1:179:A:VAL:HG21	1:182:A:SER:HA	6	0.12
(1,2491)	1:179:A:VAL:HG22	1:182:A:SER:HA	6	0.12
(1,2491)	1:179:A:VAL:HG23	1:182:A:SER:HA	6	0.12
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG21	3	0.12
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG22	3	0.12
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG23	3	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG21	3	0.12
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG22	3	0.12
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG23	3	0.12
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG21	10	0.12
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG22	10	0.12
(1,2466)	1:169:A:GLN:HG2	1:170:A:THR:HG23	10	0.12
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG21	10	0.12
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG22	10	0.12
(1,2466)	1:169:A:GLN:HG3	1:170:A:THR:HG23	10	0.12
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	5	0.12
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	5	0.12
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	13	0.12
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	13	0.12
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	13	0.12
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	13	0.12
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	13	0.12
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	13	0.12
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	15	0.12
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	15	0.12
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	15	0.12
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	15	0.12
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	15	0.12
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	15	0.12
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	16	0.12
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	16	0.12
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	16	0.12
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	16	0.12
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	16	0.12
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	16	0.12
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	19	0.12
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	19	0.12
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	19	0.12
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	19	0.12
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	19	0.12
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	19	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG11	14	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG12	14	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG13	14	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG21	14	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG22	14	0.12
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG23	14	0.12
(1,2293)	1:149:A:ARG:HD2	1:188:A:ASP:H	4	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2293)	1:149:A:ARG:HD3	1:188:A:ASP:H	4	0.12
(1,2289)	1:149:A:ARG:HG2	1:188:A:ASP:H	12	0.12
(1,2289)	1:149:A:ARG:HG3	1:188:A:ASP:H	12	0.12
(1,2267)	1:145:A:LYS:HD2	1:146:A:GLY:HA2	4	0.12
(1,2267)	1:145:A:LYS:HD3	1:146:A:GLY:HA2	4	0.12
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD11	1	0.12
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD12	1	0.12
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD13	1	0.12
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD11	1	0.12
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD12	1	0.12
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD13	1	0.12
(1,2242)	1:138:A:VAL:HG11	1:140:A:GLY:HA2	19	0.12
(1,2242)	1:138:A:VAL:HG11	1:140:A:GLY:HA3	19	0.12
(1,2242)	1:138:A:VAL:HG12	1:140:A:GLY:HA2	19	0.12
(1,2242)	1:138:A:VAL:HG12	1:140:A:GLY:HA3	19	0.12
(1,2242)	1:138:A:VAL:HG13	1:140:A:GLY:HA2	19	0.12
(1,2242)	1:138:A:VAL:HG13	1:140:A:GLY:HA3	19	0.12
(1,2242)	1:138:A:VAL:HG21	1:140:A:GLY:HA2	19	0.12
(1,2242)	1:138:A:VAL:HG21	1:140:A:GLY:HA3	19	0.12
(1,2242)	1:138:A:VAL:HG22	1:140:A:GLY:HA2	19	0.12
(1,2242)	1:138:A:VAL:HG22	1:140:A:GLY:HA3	19	0.12
(1,2242)	1:138:A:VAL:HG23	1:140:A:GLY:HA2	19	0.12
(1,2242)	1:138:A:VAL:HG23	1:140:A:GLY:HA3	19	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG11	20	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG12	20	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG13	20	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG21	20	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG22	20	0.12
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG23	20	0.12
(1,2072)	1:111:A:LEU:HD11	1:112:A:PHE:H	7	0.12
(1,2072)	1:111:A:LEU:HD12	1:112:A:PHE:H	7	0.12
(1,2072)	1:111:A:LEU:HD13	1:112:A:PHE:H	7	0.12
(1,2072)	1:111:A:LEU:HD21	1:112:A:PHE:H	7	0.12
(1,2072)	1:111:A:LEU:HD22	1:112:A:PHE:H	7	0.12
(1,2072)	1:111:A:LEU:HD23	1:112:A:PHE:H	7	0.12
(1,2044)	1:106:A:TYR:HB2	1:236:A:LYS:H	7	0.12
(1,2044)	1:106:A:TYR:HB3	1:236:A:LYS:H	7	0.12
(1,2037)	1:105:A:GLU:HG2	1:236:A:LYS:HE2	2	0.12
(1,2037)	1:105:A:GLU:HG2	1:236:A:LYS:HE3	2	0.12
(1,2037)	1:105:A:GLU:HG3	1:236:A:LYS:HE2	2	0.12
(1,2037)	1:105:A:GLU:HG3	1:236:A:LYS:HE3	2	0.12
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	5	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	5	0.12
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	5	0.12
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	5	0.12
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	5	0.12
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	5	0.12
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	7	0.12
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	7	0.12
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	7	0.12
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	7	0.12
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	7	0.12
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	7	0.12
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	12	0.12
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	12	0.12
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	12	0.12
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	12	0.12
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	12	0.12
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	12	0.12
(1,1871)	1:236:A:LYS:HG2	1:237:A:GLU:H	3	0.12
(1,1871)	1:236:A:LYS:HG3	1:237:A:GLU:H	3	0.12
(1,1871)	1:236:A:LYS:HG2	1:237:A:GLU:H	17	0.12
(1,1871)	1:236:A:LYS:HG3	1:237:A:GLU:H	17	0.12
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	2	0.12
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	2	0.12
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	2	0.12
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	2	0.12
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	2	0.12
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	2	0.12
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	2	0.12
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	2	0.12
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	2	0.12
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG21	18	0.12
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG22	18	0.12
(1,1849)	1:151:A:LEU:HD21	1:187:A:THR:HG23	18	0.12
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG21	18	0.12
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG22	18	0.12
(1,1849)	1:151:A:LEU:HD22	1:187:A:THR:HG23	18	0.12
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG21	18	0.12
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG22	18	0.12
(1,1849)	1:151:A:LEU:HD23	1:187:A:THR:HG23	18	0.12
(1,1841)	1:206:A:LYS:HD3	1:233:A:CYS:H	5	0.12
(1,1816)	1:108:A:LEU:H	1:108:A:LEU:HG	4	0.12
(1,1816)	1:108:A:LEU:H	1:108:A:LEU:HG	10	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD1	17	0.12
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD2	17	0.12
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD1	20	0.12
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD2	20	0.12
(1,1722)	1:104:A:GLU:HB2	1:236:A:LYS:HE2	8	0.12
(1,1722)	1:104:A:GLU:HB2	1:236:A:LYS:HE3	8	0.12
(1,1722)	1:104:A:GLU:HB3	1:236:A:LYS:HE2	8	0.12
(1,1722)	1:104:A:GLU:HB3	1:236:A:LYS:HE3	8	0.12
(1,1662)	1:234:A:GLN:HA	1:234:A:GLN:HG2	15	0.12
(1,1610)	1:156:A:GLU:H	1:156:A:GLU:HG3	14	0.12
(1,1337)	1:143:A:SER:HA	1:194:A:ILE:HB	4	0.12
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	1	0.12
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	11	0.12
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	14	0.12
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	20	0.12
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	9	0.12
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	9	0.12
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	9	0.12
(1,1200)	1:142:A:THR:HB	1:196:A:LEU:HB2	19	0.12
(1,1200)	1:142:A:THR:HB	1:196:A:LEU:HB3	19	0.12
(1,1150)	1:124:A:GLN:HE22	1:125:A:LEU:HD11	19	0.12
(1,1150)	1:124:A:GLN:HE22	1:125:A:LEU:HD12	19	0.12
(1,1150)	1:124:A:GLN:HE22	1:125:A:LEU:HD13	19	0.12
(1,1150)	1:124:A:GLN:HE22	1:125:A:LEU:HD21	19	0.12
(1,1150)	1:124:A:GLN:HE22	1:125:A:LEU:HD22	19	0.12
(1,1150)	1:124:A:GLN:HE22	1:125:A:LEU:HD23	19	0.12
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	2	0.12
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD2	1	0.12
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD3	1	0.12
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD2	16	0.12
(1,1110)	1:201:A:GLN:HA	1:202:A:LYS:HD3	16	0.12
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	10	0.12
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	18	0.12
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	8	0.12
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	11	0.12
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	12	0.12
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	17	0.12
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	20	0.12
(1,842)	1:133:A:GLU:H	1:146:A:GLY:HA3	10	0.12
(1,823)	1:225:A:ASN:H	1:226:A:GLY:H	11	0.12
(1,823)	1:225:A:ASN:H	1:226:A:GLY:H	13	0.12
(1,526)	1:107:A:VAL:H	1:235:A:LYS:HA	20	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,453)	1:218:A:GLY:H	1:219:A:THR:H	7	0.12
(1,453)	1:218:A:GLY:H	1:219:A:THR:H	12	0.12
(1,453)	1:218:A:GLY:H	1:219:A:THR:H	15	0.12
(1,383)	1:113:A:ASP:H	1:229:A:TYR:HA	7	0.12
(1,371)	1:194:A:ILE:HA	1:195:A:SER:H	4	0.12
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	3	0.12
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	13	0.12
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	17	0.12
(1,315)	1:104:A:GLU:HA	1:105:A:GLU:H	10	0.12
(1,315)	1:104:A:GLU:HA	1:105:A:GLU:H	20	0.12
(1,311)	1:156:A:GLU:H	1:157:A:LYS:H	15	0.12
(1,236)	1:172:A:TYR:HA	1:173:A:ASP:H	15	0.12
(1,95)	1:132:A:LEU:HD21	1:133:A:GLU:H	3	0.12
(1,95)	1:132:A:LEU:HD22	1:133:A:GLU:H	3	0.12
(1,95)	1:132:A:LEU:HD23	1:133:A:GLU:H	3	0.12
(1,2587)	1:205:A:SER:HA	1:206:A:LYS:HD2	8	0.11
(1,2587)	1:205:A:SER:HA	1:206:A:LYS:HD3	8	0.11
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG11	12	0.11
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG12	12	0.11
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG13	12	0.11
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG21	12	0.11
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG22	12	0.11
(1,2583)	1:204:A:GLY:HA2	1:232:A:VAL:HG23	12	0.11
(1,2580)	1:202:A:LYS:HB2	1:202:A:LYS:HE2	1	0.11
(1,2580)	1:202:A:LYS:HB2	1:202:A:LYS:HE3	1	0.11
(1,2580)	1:202:A:LYS:HB3	1:202:A:LYS:HE2	1	0.11
(1,2580)	1:202:A:LYS:HB3	1:202:A:LYS:HE3	1	0.11
(1,2580)	1:202:A:LYS:HB2	1:202:A:LYS:HE2	2	0.11
(1,2580)	1:202:A:LYS:HB2	1:202:A:LYS:HE3	2	0.11
(1,2580)	1:202:A:LYS:HB3	1:202:A:LYS:HE2	2	0.11
(1,2580)	1:202:A:LYS:HB3	1:202:A:LYS:HE3	2	0.11
(1,2548)	1:193:A:LYS:HG2	1:194:A:ILE:HA	13	0.11
(1,2548)	1:193:A:LYS:HG3	1:194:A:ILE:HA	13	0.11
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE21	19	0.11
(1,2460)	1:169:A:GLN:HA	1:169:A:GLN:HE22	19	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD11	2	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD12	2	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD13	2	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD21	2	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD22	2	0.11
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD23	2	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD11	2	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD12	2	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD13	2	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD21	2	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD22	2	0.11
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD23	2	0.11
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	6	0.11
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	6	0.11
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	6	0.11
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	6	0.11
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	6	0.11
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	6	0.11
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	12	0.11
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	12	0.11
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	12	0.11
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	12	0.11
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	12	0.11
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	12	0.11
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG11	18	0.11
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG12	18	0.11
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG13	18	0.11
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG21	18	0.11
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG22	18	0.11
(1,2332)	1:154:A:SER:HB2	1:217:A:VAL:HG23	18	0.11
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD11	4	0.11
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD12	4	0.11
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD13	4	0.11
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD11	4	0.11
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD12	4	0.11
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD13	4	0.11
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD11	10	0.11
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD12	10	0.11
(1,2256)	1:144:A:MET:HB2	1:194:A:ILE:HD13	10	0.11
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD11	10	0.11
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD12	10	0.11
(1,2256)	1:144:A:MET:HB3	1:194:A:ILE:HD13	10	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG11	14	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG12	14	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG13	14	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG21	14	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG22	14	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG23	14	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG11	19	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG12	19	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG13	19	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG21	19	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG22	19	0.11
(1,2155)	1:125:A:LEU:HG	1:150:A:VAL:HG23	19	0.11
(1,2123)	1:121:A:LEU:HD11	1:125:A:LEU:HA	9	0.11
(1,2123)	1:121:A:LEU:HD12	1:125:A:LEU:HA	9	0.11
(1,2123)	1:121:A:LEU:HD13	1:125:A:LEU:HA	9	0.11
(1,2123)	1:121:A:LEU:HD21	1:125:A:LEU:HA	9	0.11
(1,2123)	1:121:A:LEU:HD22	1:125:A:LEU:HA	9	0.11
(1,2123)	1:121:A:LEU:HD23	1:125:A:LEU:HA	9	0.11
(1,2088)	1:113:A:ASP:HB2	1:114:A:LEU:H	4	0.11
(1,2088)	1:113:A:ASP:HB3	1:114:A:LEU:H	4	0.11
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD11	10	0.11
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD12	10	0.11
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD13	10	0.11
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD21	10	0.11
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD22	10	0.11
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD23	10	0.11
(1,2072)	1:111:A:LEU:HD11	1:112:A:PHE:H	16	0.11
(1,2072)	1:111:A:LEU:HD12	1:112:A:PHE:H	16	0.11
(1,2072)	1:111:A:LEU:HD13	1:112:A:PHE:H	16	0.11
(1,2072)	1:111:A:LEU:HD21	1:112:A:PHE:H	16	0.11
(1,2072)	1:111:A:LEU:HD22	1:112:A:PHE:H	16	0.11
(1,2072)	1:111:A:LEU:HD23	1:112:A:PHE:H	16	0.11
(1,2072)	1:111:A:LEU:HD11	1:112:A:PHE:H	17	0.11
(1,2072)	1:111:A:LEU:HD12	1:112:A:PHE:H	17	0.11
(1,2072)	1:111:A:LEU:HD13	1:112:A:PHE:H	17	0.11
(1,2072)	1:111:A:LEU:HD21	1:112:A:PHE:H	17	0.11
(1,2072)	1:111:A:LEU:HD22	1:112:A:PHE:H	17	0.11
(1,2072)	1:111:A:LEU:HD23	1:112:A:PHE:H	17	0.11
(1,2037)	1:105:A:GLU:HG2	1:236:A:LYS:HE2	9	0.11
(1,2037)	1:105:A:GLU:HG2	1:236:A:LYS:HE3	9	0.11
(1,2037)	1:105:A:GLU:HG3	1:236:A:LYS:HE2	9	0.11
(1,2037)	1:105:A:GLU:HG3	1:236:A:LYS:HE3	9	0.11
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	13	0.11
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	13	0.11
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	13	0.11
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	13	0.11
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	13	0.11
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	13	0.11
(1,1919)	1:197:A:VAL:HG11	1:233:A:CYS:H	16	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1919)	1:197:A:VAL:HG12	1:233:A:CYS:H	16	0.11
(1,1919)	1:197:A:VAL:HG13	1:233:A:CYS:H	16	0.11
(1,1919)	1:197:A:VAL:HG21	1:233:A:CYS:H	16	0.11
(1,1919)	1:197:A:VAL:HG22	1:233:A:CYS:H	16	0.11
(1,1919)	1:197:A:VAL:HG23	1:233:A:CYS:H	16	0.11
(1,1902)	1:103:A:THR:H	1:103:A:THR:HG21	17	0.11
(1,1902)	1:103:A:THR:H	1:103:A:THR:HG22	17	0.11
(1,1902)	1:103:A:THR:H	1:103:A:THR:HG23	17	0.11
(1,1874)	1:165:A:LEU:HD11	1:208:A:GLU:H	5	0.11
(1,1874)	1:165:A:LEU:HD12	1:208:A:GLU:H	5	0.11
(1,1874)	1:165:A:LEU:HD13	1:208:A:GLU:H	5	0.11
(1,1871)	1:236:A:LYS:HG2	1:237:A:GLU:H	19	0.11
(1,1871)	1:236:A:LYS:HG3	1:237:A:GLU:H	19	0.11
(1,1841)	1:206:A:LYS:HD3	1:233:A:CYS:H	3	0.11
(1,1816)	1:108:A:LEU:H	1:108:A:LEU:HG	3	0.11
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD1	3	0.11
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD2	3	0.11
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD1	11	0.11
(1,1795)	1:132:A:LEU:HG	1:229:A:TYR:HD2	11	0.11
(1,1767)	1:120:A:ASP:HA	1:123:A:GLN:HB2	6	0.11
(1,1767)	1:120:A:ASP:HA	1:123:A:GLN:HB3	6	0.11
(1,1765)	1:160:A:TYR:HD1	1:212:A:ARG:HG3	5	0.11
(1,1765)	1:160:A:TYR:HD2	1:212:A:ARG:HG3	5	0.11
(1,1765)	1:160:A:TYR:HD1	1:212:A:ARG:HG3	17	0.11
(1,1765)	1:160:A:TYR:HD2	1:212:A:ARG:HG3	17	0.11
(1,1747)	1:182:A:SER:H	1:183:A:CYS:HB2	5	0.11
(1,1637)	1:206:A:LYS:HB2	1:231:A:LEU:H	9	0.11
(1,1637)	1:206:A:LYS:HB3	1:231:A:LEU:H	9	0.11
(1,1610)	1:156:A:GLU:H	1:156:A:GLU:HG3	15	0.11
(1,1577)	1:147:A:ILE:HB	1:191:A:SER:HB2	6	0.11
(1,1405)	1:137:A:TYR:HE1	1:234:A:GLN:HA	1	0.11
(1,1405)	1:137:A:TYR:HE2	1:234:A:GLN:HA	1	0.11
(1,1405)	1:137:A:TYR:HE1	1:234:A:GLN:HA	8	0.11
(1,1405)	1:137:A:TYR:HE2	1:234:A:GLN:HA	8	0.11
(1,1350)	1:172:A:TYR:HE1	1:209:A:PHE:HA	11	0.11
(1,1350)	1:172:A:TYR:HE2	1:209:A:PHE:HA	11	0.11
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	3	0.11
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	7	0.11
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	10	0.11
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	15	0.11
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	16	0.11
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	18	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1273)	1:172:A:TYR:HE1	1:174:A:ILE:HA	5	0.11
(1,1273)	1:172:A:TYR:HE2	1:174:A:ILE:HA	5	0.11
(1,1273)	1:172:A:TYR:HE1	1:174:A:ILE:HA	12	0.11
(1,1273)	1:172:A:TYR:HE2	1:174:A:ILE:HA	12	0.11
(1,1246)	1:205:A:SER:HB2	1:206:A:LYS:HE2	19	0.11
(1,1246)	1:205:A:SER:HB2	1:206:A:LYS:HE3	19	0.11
(1,1246)	1:205:A:SER:HB3	1:206:A:LYS:HE2	19	0.11
(1,1246)	1:205:A:SER:HB3	1:206:A:LYS:HE3	19	0.11
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	10	0.11
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	10	0.11
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	10	0.11
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	14	0.11
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	14	0.11
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	14	0.11
(1,1224)	1:153:A:ILE:HD11	1:154:A:SER:HB3	19	0.11
(1,1224)	1:153:A:ILE:HD12	1:154:A:SER:HB3	19	0.11
(1,1224)	1:153:A:ILE:HD13	1:154:A:SER:HB3	19	0.11
(1,1191)	1:224:A:ASN:HD21	1:227:A:THR:HB	12	0.11
(1,1191)	1:224:A:ASN:HD22	1:227:A:THR:HB	12	0.11
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	8	0.11
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	9	0.11
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	12	0.11
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG21	15	0.11
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG22	15	0.11
(1,1146)	1:112:A:PHE:HD1	1:135:A:THR:HG23	15	0.11
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG21	15	0.11
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG22	15	0.11
(1,1146)	1:112:A:PHE:HD2	1:135:A:THR:HG23	15	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD11	2	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD12	2	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD13	2	0.11
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD11	2	0.11
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD12	2	0.11
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD13	2	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD11	2	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD12	2	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD13	2	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD11	5	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD12	5	0.11
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD13	5	0.11
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD11	5	0.11
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD12	5	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD13	5	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD11	5	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD12	5	0.11
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD13	5	0.11
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG21	2	0.11
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG22	2	0.11
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG23	2	0.11
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG21	2	0.11
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG22	2	0.11
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG23	2	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	1	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	2	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	3	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	4	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	5	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	6	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	7	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	8	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	9	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	11	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	12	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	13	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	14	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	15	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	16	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	17	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	19	0.11
(1,1097)	1:127:A:VAL:HA	1:127:A:VAL:HB	20	0.11
(1,1088)	1:236:A:LYS:HE2	1:237:A:GLU:HA	6	0.11
(1,1088)	1:236:A:LYS:HE3	1:237:A:GLU:HA	6	0.11
(1,1088)	1:236:A:LYS:HE2	1:237:A:GLU:HA	9	0.11
(1,1088)	1:236:A:LYS:HE3	1:237:A:GLU:HA	9	0.11
(1,1062)	1:114:A:LEU:HD21	1:115:A:PRO:HD3	1	0.11
(1,1062)	1:114:A:LEU:HD22	1:115:A:PRO:HD3	1	0.11
(1,1062)	1:114:A:LEU:HD23	1:115:A:PRO:HD3	1	0.11
(1,1062)	1:114:A:LEU:HD21	1:115:A:PRO:HD3	16	0.11
(1,1062)	1:114:A:LEU:HD22	1:115:A:PRO:HD3	16	0.11
(1,1062)	1:114:A:LEU:HD23	1:115:A:PRO:HD3	16	0.11
(1,1062)	1:114:A:LEU:HD21	1:115:A:PRO:HD3	19	0.11
(1,1062)	1:114:A:LEU:HD22	1:115:A:PRO:HD3	19	0.11
(1,1062)	1:114:A:LEU:HD23	1:115:A:PRO:HD3	19	0.11
(1,1055)	1:154:A:SER:HB3	1:157:A:LYS:HB3	1	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1055)	1:154:A:SER:HB3	1:157:A:LYS:HB3	18	0.11
(1,968)	1:152:A:ASN:HD22	1:153:A:ILE:HA	5	0.11
(1,968)	1:152:A:ASN:HD22	1:153:A:ILE:HA	6	0.11
(1,943)	1:128:A:GLN:HE21	1:129:A:LYS:HD2	7	0.11
(1,894)	1:142:A:THR:H	1:196:A:LEU:HA	17	0.11
(1,757)	1:125:A:LEU:HA	1:126:A:GLN:HE21	9	0.11
(1,655)	1:126:A:GLN:H	1:129:A:LYS:H	10	0.11
(1,581)	1:128:A:GLN:H	1:128:A:GLN:HG3	8	0.11
(1,526)	1:107:A:VAL:H	1:235:A:LYS:HA	13	0.11
(1,453)	1:218:A:GLY:H	1:219:A:THR:H	19	0.11
(1,371)	1:194:A:ILE:HA	1:195:A:SER:H	15	0.11
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	8	0.11
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	16	0.11
(1,362)	1:135:A:THR:HB	1:136:A:GLU:H	13	0.11
(1,268)	1:104:A:GLU:H	1:104:A:GLU:HG2	16	0.11
(1,268)	1:104:A:GLU:H	1:104:A:GLU:HG3	16	0.11
(1,238)	1:173:A:ASP:H	1:173:A:ASP:HB3	4	0.11
(1,113)	1:161:A:VAL:HG21	1:212:A:ARG:H	1	0.11
(1,113)	1:161:A:VAL:HG22	1:212:A:ARG:H	1	0.11
(1,113)	1:161:A:VAL:HG23	1:212:A:ARG:H	1	0.11
(1,113)	1:161:A:VAL:HG21	1:212:A:ARG:H	10	0.11
(1,113)	1:161:A:VAL:HG22	1:212:A:ARG:H	10	0.11
(1,113)	1:161:A:VAL:HG23	1:212:A:ARG:H	10	0.11
(1,2622)	1:210:A:CYS:HB2	1:223:A:ASN:HB2	16	0.1
(1,2622)	1:210:A:CYS:HB2	1:223:A:ASN:HB3	16	0.1
(1,2622)	1:210:A:CYS:HB3	1:223:A:ASN:HB2	16	0.1
(1,2622)	1:210:A:CYS:HB3	1:223:A:ASN:HB3	16	0.1
(1,2579)	1:201:A:GLN:HE21	1:202:A:LYS:H	11	0.1
(1,2579)	1:201:A:GLN:HE22	1:202:A:LYS:H	11	0.1
(1,2485)	1:179:A:VAL:HB	1:180:A:PRO:HD2	1	0.1
(1,2485)	1:179:A:VAL:HB	1:180:A:PRO:HD3	1	0.1
(1,2462)	1:169:A:GLN:HB2	1:170:A:THR:H	9	0.1
(1,2462)	1:169:A:GLN:HB3	1:170:A:THR:H	9	0.1
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD11	3	0.1
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD12	3	0.1
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD13	3	0.1
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD21	3	0.1
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD22	3	0.1
(1,2390)	1:160:A:TYR:HD1	1:175:A:LEU:HD23	3	0.1
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD11	3	0.1
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD12	3	0.1
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD13	3	0.1

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD21	3	0.1
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD22	3	0.1
(1,2390)	1:160:A:TYR:HD2	1:175:A:LEU:HD23	3	0.1
(1,2368)	1:159:A:VAL:HG11	1:160:A:TYR:H	17	0.1
(1,2368)	1:159:A:VAL:HG12	1:160:A:TYR:H	17	0.1
(1,2368)	1:159:A:VAL:HG13	1:160:A:TYR:H	17	0.1
(1,2368)	1:159:A:VAL:HG21	1:160:A:TYR:H	17	0.1
(1,2368)	1:159:A:VAL:HG22	1:160:A:TYR:H	17	0.1
(1,2368)	1:159:A:VAL:HG23	1:160:A:TYR:H	17	0.1
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	3	0.1
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	3	0.1
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	3	0.1
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	3	0.1
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	3	0.1
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	3	0.1
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	4	0.1
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	4	0.1
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	4	0.1
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	4	0.1
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	4	0.1
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	4	0.1
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	10	0.1
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	10	0.1
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	10	0.1
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	10	0.1
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	10	0.1
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	10	0.1
(1,2362)	1:158:A:LEU:HD11	1:160:A:TYR:HA	20	0.1
(1,2362)	1:158:A:LEU:HD12	1:160:A:TYR:HA	20	0.1
(1,2362)	1:158:A:LEU:HD13	1:160:A:TYR:HA	20	0.1
(1,2362)	1:158:A:LEU:HD21	1:160:A:TYR:HA	20	0.1
(1,2362)	1:158:A:LEU:HD22	1:160:A:TYR:HA	20	0.1
(1,2362)	1:158:A:LEU:HD23	1:160:A:TYR:HA	20	0.1
(1,2297)	1:150:A:VAL:H	1:189:A:GLN:HG2	11	0.1
(1,2297)	1:150:A:VAL:H	1:189:A:GLN:HG3	11	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD11	3	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD12	3	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD13	3	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD21	3	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD22	3	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD23	3	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD11	20	0.1

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD12	20	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD13	20	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD21	20	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD22	20	0.1
(1,2079)	1:112:A:PHE:H	1:231:A:LEU:HD23	20	0.1
(1,2044)	1:106:A:TYR:HB2	1:236:A:LYS:H	5	0.1
(1,2044)	1:106:A:TYR:HB3	1:236:A:LYS:H	5	0.1
(1,1854)	1:108:A:LEU:HD21	1:205:A:SER:HB2	15	0.1
(1,1854)	1:108:A:LEU:HD21	1:205:A:SER:HB3	15	0.1
(1,1854)	1:108:A:LEU:HD22	1:205:A:SER:HB2	15	0.1
(1,1854)	1:108:A:LEU:HD22	1:205:A:SER:HB3	15	0.1
(1,1854)	1:108:A:LEU:HD23	1:205:A:SER:HB2	15	0.1
(1,1854)	1:108:A:LEU:HD23	1:205:A:SER:HB3	15	0.1
(1,1816)	1:108:A:LEU:H	1:108:A:LEU:HG	7	0.1
(1,1816)	1:108:A:LEU:H	1:108:A:LEU:HG	14	0.1
(1,1747)	1:182:A:SER:H	1:183:A:CYS:HB2	12	0.1
(1,1577)	1:147:A:ILE:HB	1:191:A:SER:HB2	11	0.1
(1,1499)	1:196:A:LEU:HB2	1:200:A:TYR:HE1	9	0.1
(1,1499)	1:196:A:LEU:HB2	1:200:A:TYR:HE2	9	0.1
(1,1499)	1:196:A:LEU:HB3	1:200:A:TYR:HE1	9	0.1
(1,1499)	1:196:A:LEU:HB3	1:200:A:TYR:HE2	9	0.1
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD1	8	0.1
(1,1487)	1:108:A:LEU:HB3	1:137:A:TYR:HD2	8	0.1
(1,1337)	1:143:A:SER:HA	1:194:A:ILE:HB	13	0.1
(1,1304)	1:197:A:VAL:HA	1:198:A:PRO:HB3	4	0.1
(1,1216)	1:143:A:SER:HB2	1:193:A:LYS:HD2	16	0.1
(1,1216)	1:143:A:SER:HB2	1:193:A:LYS:HD3	16	0.1
(1,1216)	1:143:A:SER:HB3	1:193:A:LYS:HD2	16	0.1
(1,1216)	1:143:A:SER:HB3	1:193:A:LYS:HD3	16	0.1
(1,1147)	1:128:A:GLN:HE21	1:129:A:LYS:HG2	5	0.1
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD11	8	0.1
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD12	8	0.1
(1,1138)	1:135:A:THR:HG21	1:194:A:ILE:HD13	8	0.1
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD11	8	0.1
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD12	8	0.1
(1,1138)	1:135:A:THR:HG22	1:194:A:ILE:HD13	8	0.1
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD11	8	0.1
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD12	8	0.1
(1,1138)	1:135:A:THR:HG23	1:194:A:ILE:HD13	8	0.1
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG21	14	0.1
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG22	14	0.1
(1,1117)	1:206:A:LYS:HE2	1:230:A:THR:HG23	14	0.1

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG21	14	0.1
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG22	14	0.1
(1,1117)	1:206:A:LYS:HE3	1:230:A:THR:HG23	14	0.1
(1,1042)	1:198:A:PRO:HB2	1:199:A:PRO:HA	16	0.1
(1,968)	1:152:A:ASN:HD22	1:153:A:ILE:HA	2	0.1
(1,665)	1:152:A:ASN:HD22	1:185:A:GLY:H	18	0.1
(1,662)	1:181:A:ASN:HA	1:181:A:ASN:HD21	12	0.1
(1,581)	1:128:A:GLN:H	1:128:A:GLN:HG3	11	0.1
(1,526)	1:107:A:VAL:H	1:235:A:LYS:HA	16	0.1
(1,453)	1:218:A:GLY:H	1:219:A:THR:H	6	0.1
(1,383)	1:113:A:ASP:H	1:229:A:TYR:HA	19	0.1
(1,371)	1:194:A:ILE:HA	1:195:A:SER:H	5	0.1
(1,366)	1:152:A:ASN:HD21	1:186:A:GLU:H	5	0.1
(1,345)	1:160:A:TYR:HB2	1:161:A:VAL:H	8	0.1
(1,345)	1:160:A:TYR:HB3	1:161:A:VAL:H	8	0.1
(1,313)	1:157:A:LYS:H	1:157:A:LYS:HB2	14	0.1
(1,269)	1:104:A:GLU:H	1:104:A:GLU:HB2	10	0.1
(1,269)	1:104:A:GLU:H	1:104:A:GLU:HB3	10	0.1
(1,238)	1:173:A:ASP:H	1:173:A:ASP:HB3	6	0.1
(1,95)	1:132:A:LEU:HD21	1:133:A:GLU:H	8	0.1
(1,95)	1:132:A:LEU:HD22	1:133:A:GLU:H	8	0.1
(1,95)	1:132:A:LEU:HD23	1:133:A:GLU:H	8	0.1

10 Dihedral-angle violation analysis ⓘ

No dihedral-angle restraints found