



Full wwPDB NMR Structure Validation Report ⓘ

Jun 5, 2023 – 06:06 PM JST

PDB ID : 6IVU
BMRB ID : 36221
Title : Solution structure of the Sigma-anti-sigma factor complex RsgI1N-SigI1C from Clostridium thermocellum
Authors : Wei, Z.; Feng, Y.
Deposited on : 2018-12-04

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 : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
BMRB Restraints Analysis : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.33

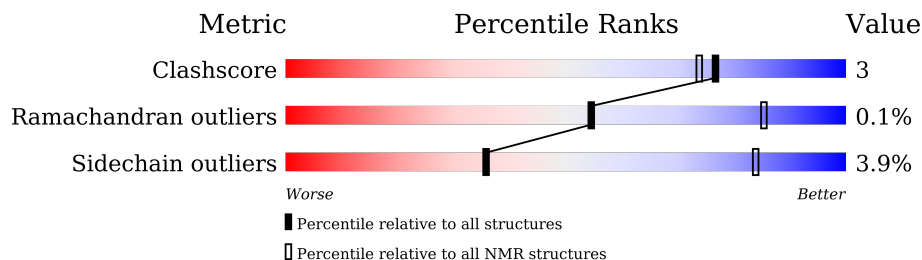
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 92%.

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



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

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

Mol	Chain	Length	Quality of chain
1	A	53	 81% . 15%
2	B	118	 83% 7% . 9%

2 Ensemble composition and analysis i

This entry contains 20 models. Model 16 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:3-A:47, B:143-B:249 (152)	0.40	16

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

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

The models can be grouped into 2 clusters. No single-model clusters were found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2909 atoms, of which 1482 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Anti-sigma-I factor RsgI1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	53	879	270	448	78	80	3	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP A3DBH1

- Molecule 2 is a protein called RNA polymerase sigma factor SigI1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
2	B	118	2030	629	1034	186	174	7	0

There are 9 discrepancies between the modelled and reference sequences:

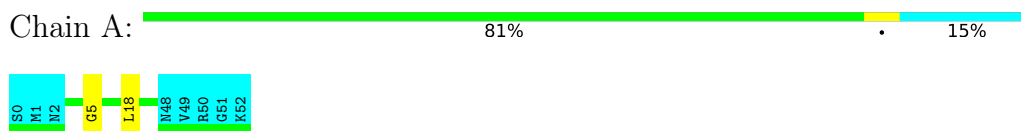
Chain	Residue	Modelled	Actual	Comment	Reference
B	140	MET	-	initiating methionine	UNP A3DBH0
B	250	LEU	-	expression tag	UNP A3DBH0
B	251	GLU	-	expression tag	UNP A3DBH0
B	252	HIS	-	expression tag	UNP A3DBH0
B	253	HIS	-	expression tag	UNP A3DBH0
B	254	HIS	-	expression tag	UNP A3DBH0
B	255	HIS	-	expression tag	UNP A3DBH0
B	256	HIS	-	expression tag	UNP A3DBH0
B	257	HIS	-	expression tag	UNP A3DBH0

4 Residue-property plots

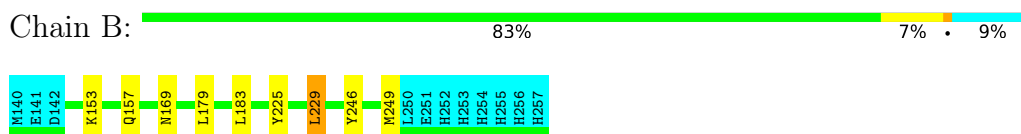
4.1 Average score per residue in the NMR ensemble

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: Anti-sigma-I factor RsgI1



- Molecule 2: RNA polymerase sigma factor SigI1

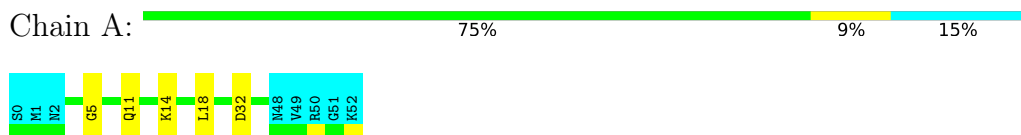


4.2 Scores per residue for each member of the ensemble

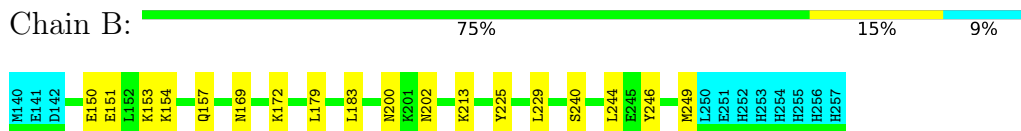
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: Anti-sigma-I factor RsgI1

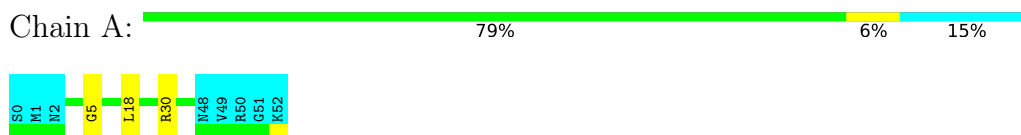


- Molecule 2: RNA polymerase sigma factor SigI1

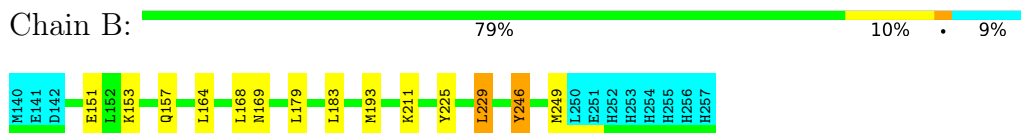


4.2.2 Score per residue for model 2

- Molecule 1: Anti-sigma-I factor RsgI1

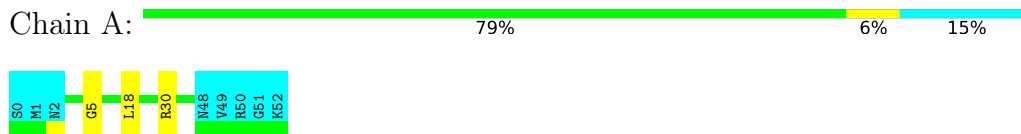


- Molecule 2: RNA polymerase sigma factor SigI1

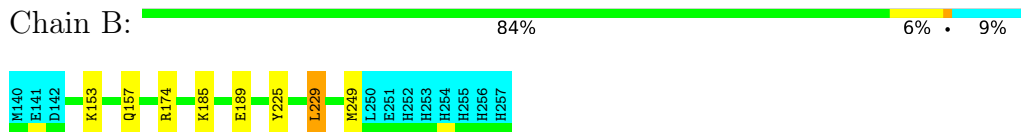


4.2.3 Score per residue for model 3

- Molecule 1: Anti-sigma-I factor RsgI1

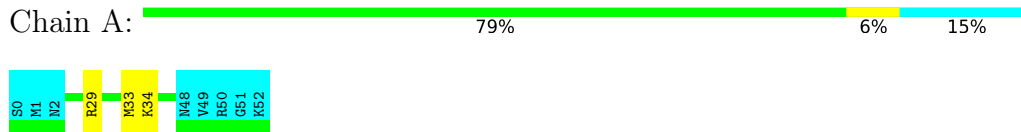


- Molecule 2: RNA polymerase sigma factor SigI1

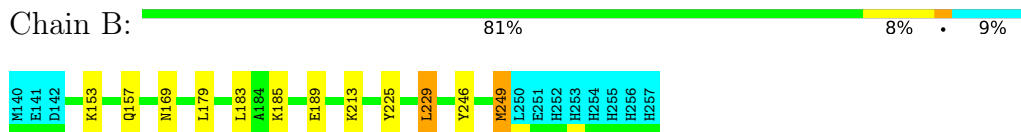


4.2.4 Score per residue for model 4

- Molecule 1: Anti-sigma-I factor RsgI1

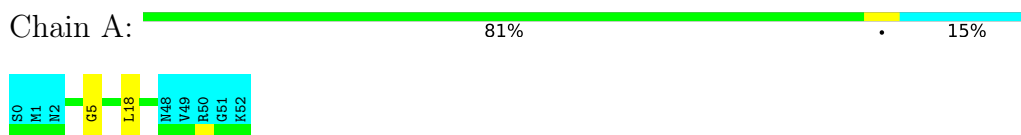


- Molecule 2: RNA polymerase sigma factor SigI1

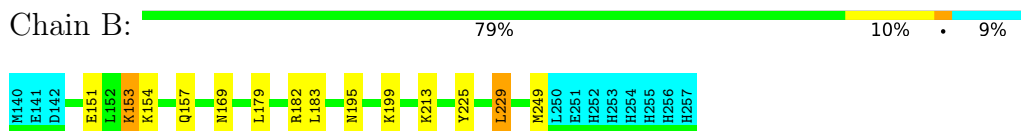


4.2.5 Score per residue for model 5

- Molecule 1: Anti-sigma-I factor RsgI1

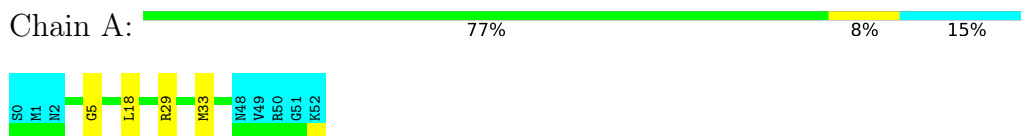


- Molecule 2: RNA polymerase sigma factor SigI1

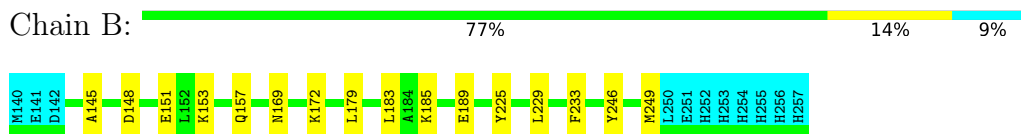


4.2.6 Score per residue for model 6

- Molecule 1: Anti-sigma-I factor RsgI1

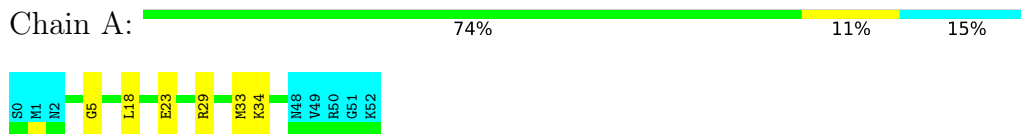


- Molecule 2: RNA polymerase sigma factor SigI1

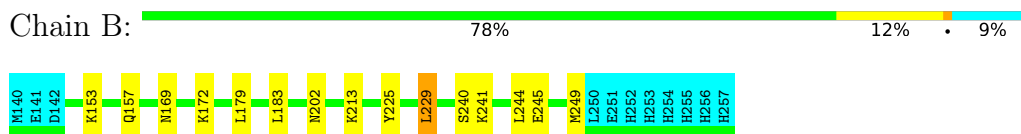


4.2.7 Score per residue for model 7

- Molecule 1: Anti-sigma-I factor RsgI1

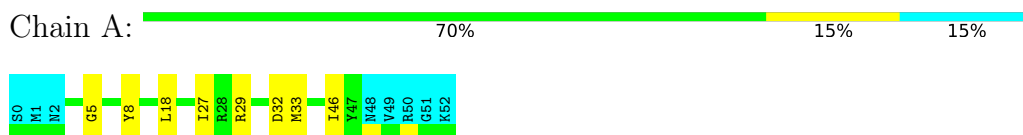


- Molecule 2: RNA polymerase sigma factor SigI1

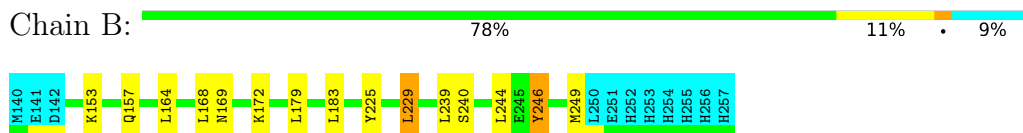


4.2.8 Score per residue for model 8

- Molecule 1: Anti-sigma-I factor RsgI1

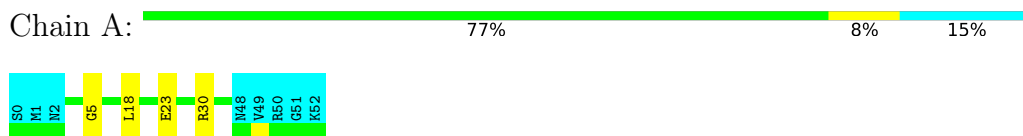


- Molecule 2: RNA polymerase sigma factor SigI1

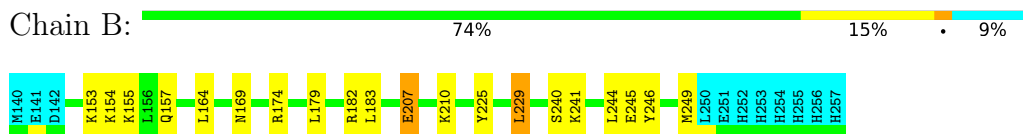


4.2.9 Score per residue for model 9

- Molecule 1: Anti-sigma-I factor RsgI1

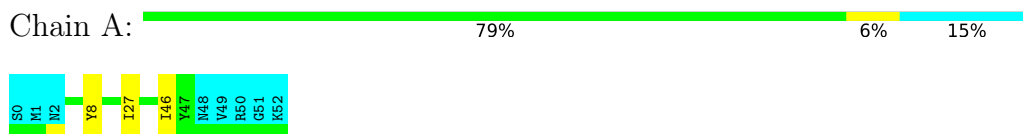


- Molecule 2: RNA polymerase sigma factor SigI1

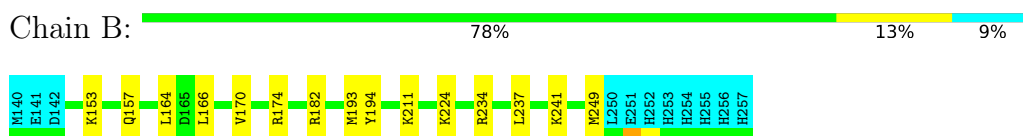


4.2.10 Score per residue for model 10

- Molecule 1: Anti-sigma-I factor RsgI1

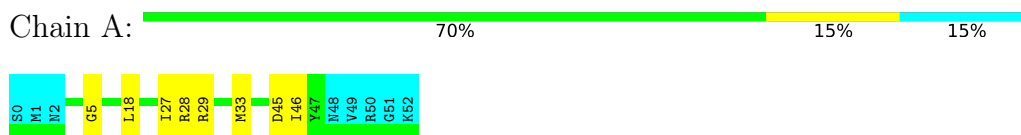


- Molecule 2: RNA polymerase sigma factor SigI1

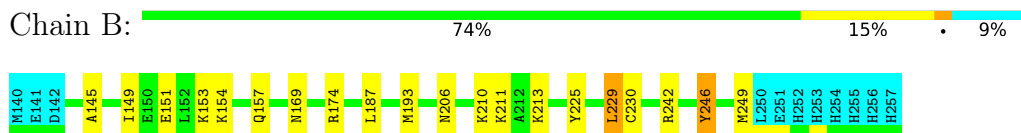


4.2.11 Score per residue for model 11

- Molecule 1: Anti-sigma-I factor RsgI1

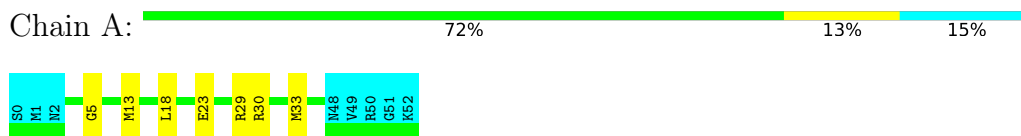


- Molecule 2: RNA polymerase sigma factor SigI1

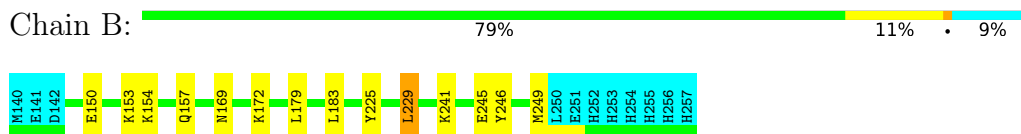


4.2.12 Score per residue for model 12

- Molecule 1: Anti-sigma-I factor RsgI1

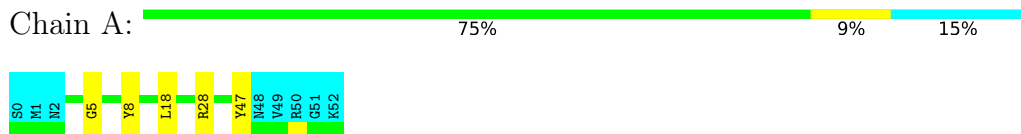


- Molecule 2: RNA polymerase sigma factor SigI1

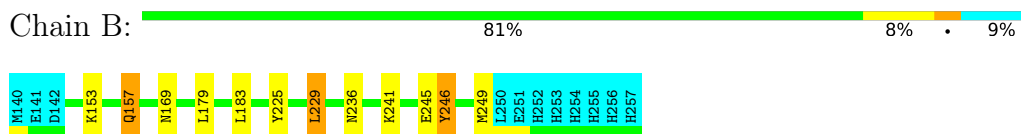


4.2.13 Score per residue for model 13

- Molecule 1: Anti-sigma-I factor RsgI1

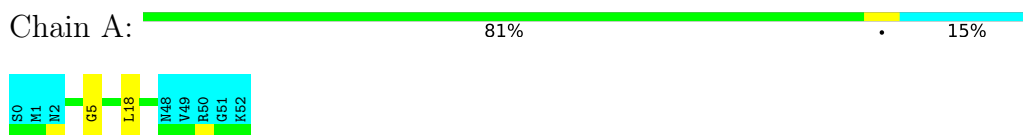


- Molecule 2: RNA polymerase sigma factor SigI1

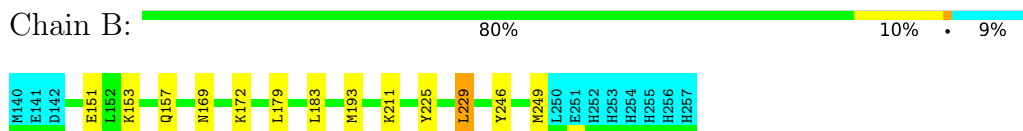


4.2.14 Score per residue for model 14

- Molecule 1: Anti-sigma-I factor RsgI1

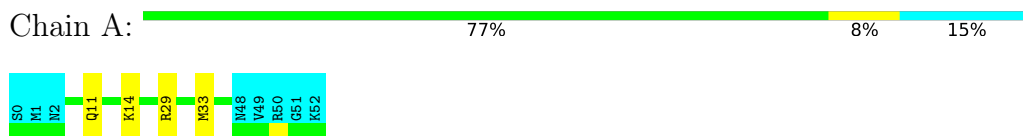


- Molecule 2: RNA polymerase sigma factor SigI1

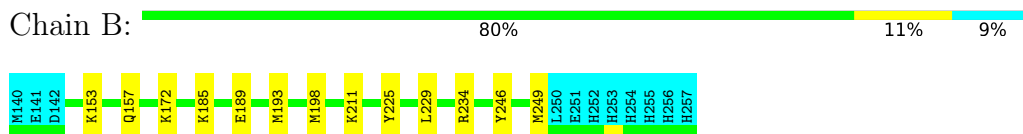


4.2.15 Score per residue for model 15

- Molecule 1: Anti-sigma-I factor RsgI1

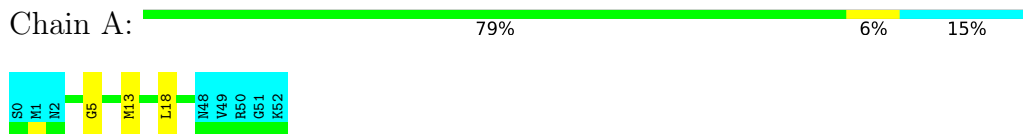


- Molecule 2: RNA polymerase sigma factor SigI1

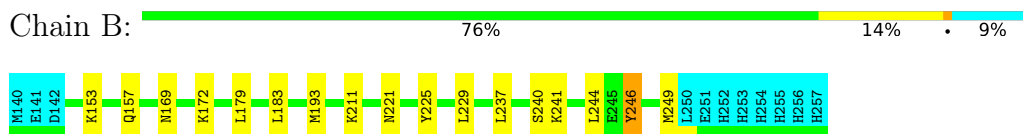


4.2.16 Score per residue for model 16 (medoid)

- Molecule 1: Anti-sigma-I factor RsgI1

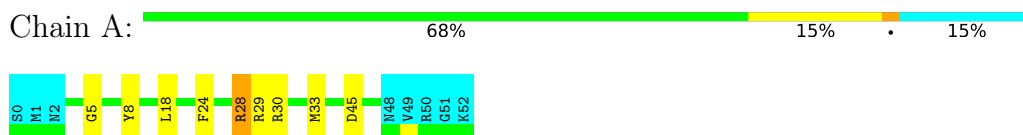


- Molecule 2: RNA polymerase sigma factor SigI1

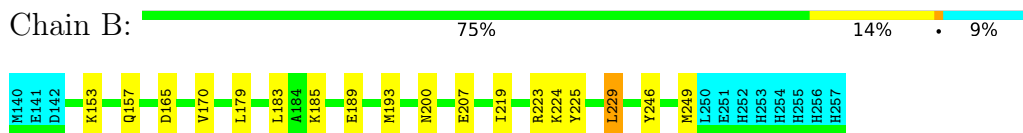


4.2.17 Score per residue for model 17

- Molecule 1: Anti-sigma-I factor RsgI1

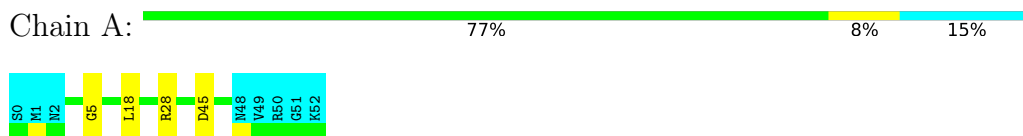


- Molecule 2: RNA polymerase sigma factor SigI1

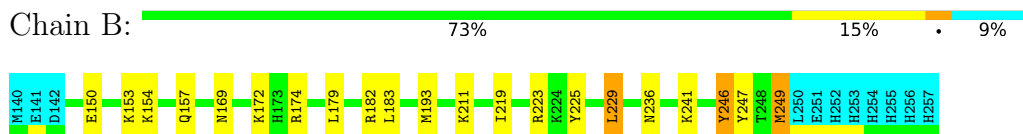


4.2.18 Score per residue for model 18

- Molecule 1: Anti-sigma-I factor RsgI1

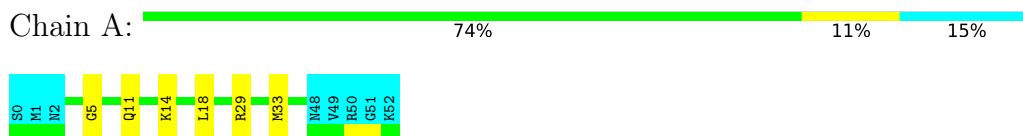


- Molecule 2: RNA polymerase sigma factor SigI1

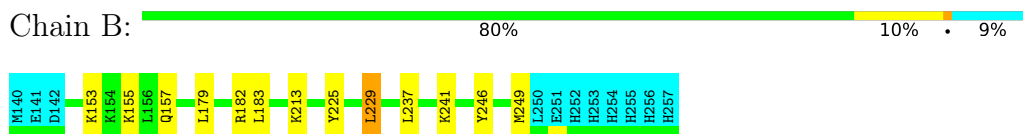


4.2.19 Score per residue for model 19

- Molecule 1: Anti-sigma-I factor RsgI1

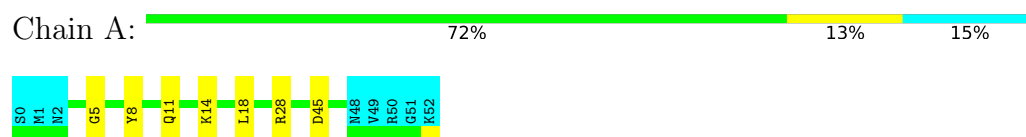


- Molecule 2: RNA polymerase sigma factor SigI1

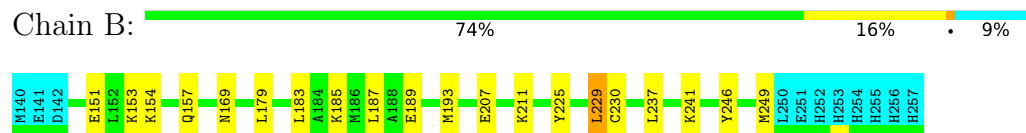


4.2.20 Score per residue for model 20

- Molecule 1: Anti-sigma-I factor RsgI1



- Molecule 2: RNA polymerase sigma factor SigI1



5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
CNS	refinement	
CNS	structure calculation	

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	2245
Number of shifts mapped to atoms	2245
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	92%

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

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	0.3±0.5
2	B	0.0±0.0	0.4±0.6
All	All	0	14

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	30	ARG	Sidechain	5
2	B	182	ARG	Sidechain	5
2	B	174	ARG	Sidechain	1
2	B	242	ARG	Sidechain	1
1	A	28	ARG	Sidechain	1
2	B	234	ARG	Sidechain	1

6.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	369	382	382	2±1
2	B	893	948	946	6±2
All	All	25240	26600	26560	153

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 3.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:150:GLU:O	2:B:154:LYS:HG3	0.59	1.98	1	3
1:A:29:ARG:HA	1:A:33:MET:SD	0.58	2.38	4	9
1:A:5:GLY:HA3	1:A:18:LEU:O	0.56	2.01	1	16
2:B:153:LYS:O	2:B:157:GLN:HG2	0.51	2.05	19	20
2:B:237:LEU:O	2:B:241:LYS:HB2	0.50	2.07	19	4
2:B:225:TYR:O	2:B:229:LEU:HD22	0.50	2.07	6	19
2:B:151:GLU:O	2:B:154:LYS:HG2	0.50	2.07	5	2
1:A:11:GLN:O	1:A:14:LYS:HG2	0.49	2.08	19	4
2:B:219:ILE:O	2:B:223:ARG:HB3	0.49	2.08	18	2
2:B:179:LEU:O	2:B:183:LEU:HG	0.49	2.08	17	16
2:B:185:LYS:O	2:B:189:GLU:HG2	0.48	2.08	15	6
2:B:193:MET:SD	2:B:207:GLU:HB3	0.48	2.49	17	1
1:A:18:LEU:HA	1:A:23:GLU:O	0.48	2.09	7	2
1:A:28:ARG:HD2	1:A:45:ASP:OD1	0.47	2.10	20	2
2:B:187:LEU:HG	2:B:230:CYS:SG	0.47	2.49	11	2
1:A:23:GLU:HA	2:B:225:TYR:OH	0.46	2.10	9	1
1:A:28:ARG:NE	1:A:45:ASP:HA	0.46	2.26	17	1
2:B:241:LYS:O	2:B:245:GLU:HG3	0.46	2.11	9	4
2:B:165:ASP:HB3	2:B:224:LYS:HD3	0.45	1.88	17	1
2:B:174:ARG:HG2	2:B:247:TYR:OH	0.45	2.12	18	1
2:B:240:SER:O	2:B:244:LEU:HG	0.44	2.12	9	5
1:A:27:ILE:HG22	1:A:46:ILE:HA	0.44	1.89	11	3
2:B:151:GLU:OE1	2:B:154:LYS:HE3	0.44	2.12	11	1
2:B:172:LYS:N	2:B:172:LYS:HD2	0.44	2.28	7	4
2:B:145:ALA:HB3	2:B:148:ASP:OD1	0.44	2.13	6	1
1:A:28:ARG:HD2	1:A:45:ASP:OD2	0.44	2.12	11	1
1:A:5:GLY:HA2	1:A:18:LEU:O	0.44	2.13	8	1
2:B:166:LEU:O	2:B:170:VAL:HG23	0.44	2.13	10	1
2:B:193:MET:SD	2:B:211:LYS:HD2	0.43	2.53	18	6
2:B:207:GLU:O	2:B:210:LYS:HG2	0.43	2.13	9	1
2:B:145:ALA:O	2:B:149:ILE:HG12	0.43	2.14	11	1
2:B:193:MET:SD	2:B:211:LYS:HD3	0.43	2.54	20	2
2:B:246:TYR:O	2:B:249:MET:HG2	0.43	2.14	2	2
2:B:164:LEU:O	2:B:168:LEU:HG	0.42	2.14	8	2
2:B:246:TYR:HA	2:B:249:MET:HG3	0.42	1.91	18	1
1:A:24:PHE:O	2:B:170:VAL:HG21	0.41	2.14	17	1
2:B:194:TYR:CE1	2:B:234:ARG:HG3	0.41	2.50	10	1
2:B:195:ASN:O	2:B:199:LYS:HG3	0.40	2.15	5	1
2:B:185:LYS:HG2	2:B:233:PHE:CZ	0.40	2.51	6	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:206:ASN:O	2:B:210:LYS:HG2	0.40	2.16	11	1

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	45/53 (85%)	43±1 (96±2%)	2±1 (4±2%)	0±0 (0±0%)	54	85
2	B	107/118 (91%)	104±1 (97±1%)	3±1 (3±1%)	0±0 (0±0%)	54	85
All	All	3040/3420 (89%)	2948 (97%)	89 (3%)	3 (0%)	54	85

All 2 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	B	236	ASN	2
1	A	47	TYR	1

6.3.2 Protein sidechains [i](#)

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	40/47 (85%)	39±1 (98±2%)	1±1 (2±2%)	66	95
2	B	98/109 (90%)	93±1 (95±1%)	5±1 (5±1%)	29	78
All	All	2760/3120 (88%)	2652 (96%)	108 (4%)	36	84

All 25 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	B	249	MET	18
2	B	169	ASN	15
2	B	229	LEU	15
2	B	246	TYR	13
2	B	213	LYS	6
2	B	172	LYS	5
1	A	8	TYR	5
2	B	151	GLU	4
2	B	174	ARG	3
1	A	32	ASP	2
2	B	200	ASN	2
2	B	202	ASN	2
1	A	34	LYS	2
2	B	155	LYS	2
2	B	164	LEU	2
2	B	207	GLU	2
1	A	13	MET	2
2	B	153	LYS	1
2	B	154	LYS	1
2	B	224	LYS	1
2	B	157	GLN	1
2	B	198	MET	1
2	B	221	ASN	1
1	A	28	ARG	1
2	B	241	LYS	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

6.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation [i](#)

The completeness of assignment taking into account all chemical shift lists is 92% for the well-defined parts and 88% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_0*

7.1.1 Bookkeeping [i](#)

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

Total number of shifts	2245
Number of shifts mapped to atoms	2245
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	2

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	164	-0.57 ± 0.13	Should be checked
$^{13}\text{C}_\beta$	157	-0.08 ± 0.10	None needed (< 0.5 ppm)
$^{13}\text{C}'$	164	-0.28 ± 0.07	None needed (< 0.5 ppm)
^{15}N	157	0.43 ± 0.19	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 92%, i.e. 2098 atoms were assigned a chemical shift out of a possible 2278. 0 out of 25 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	753/763 (99%)	304/309 (98%)	302/304 (99%)	147/150 (98%)
Sidechain	1235/1386 (89%)	838/894 (94%)	372/421 (88%)	25/71 (35%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	110/129 (85%)	55/61 (90%)	55/64 (86%)	0/4 (0%)
Overall	2098/2278 (92%)	1197/1264 (95%)	729/789 (92%)	172/225 (76%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 88%, i.e. 2245 atoms were assigned a chemical shift out of a possible 2549. 0 out of 27 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	813/859 (95%)	328/348 (94%)	328/342 (96%)	157/169 (93%)
Sidechain	1322/1513 (87%)	897/975 (92%)	398/461 (86%)	27/77 (35%)
Aromatic	110/177 (62%)	55/85 (65%)	55/76 (72%)	0/16 (0%)
Overall	2245/2549 (88%)	1280/1408 (91%)	781/879 (89%)	184/262 (70%)

7.1.4 Statistically unusual chemical shifts [i](#)

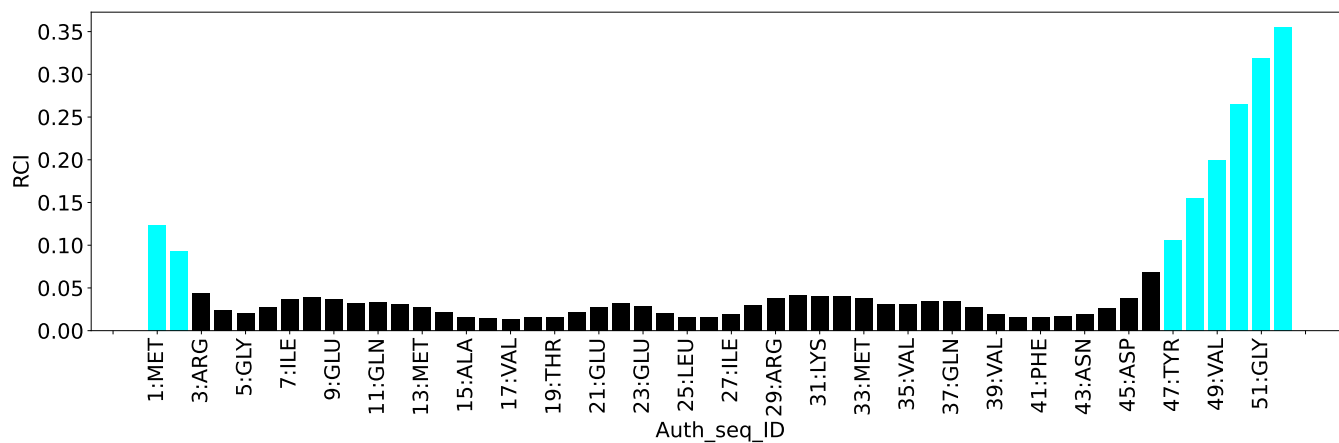
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	3	ARG	HB2	0.28	0.52 – 3.08	-5.9
1	B	201	LYS	HD3	0.54	0.54 – 2.65	-5.0

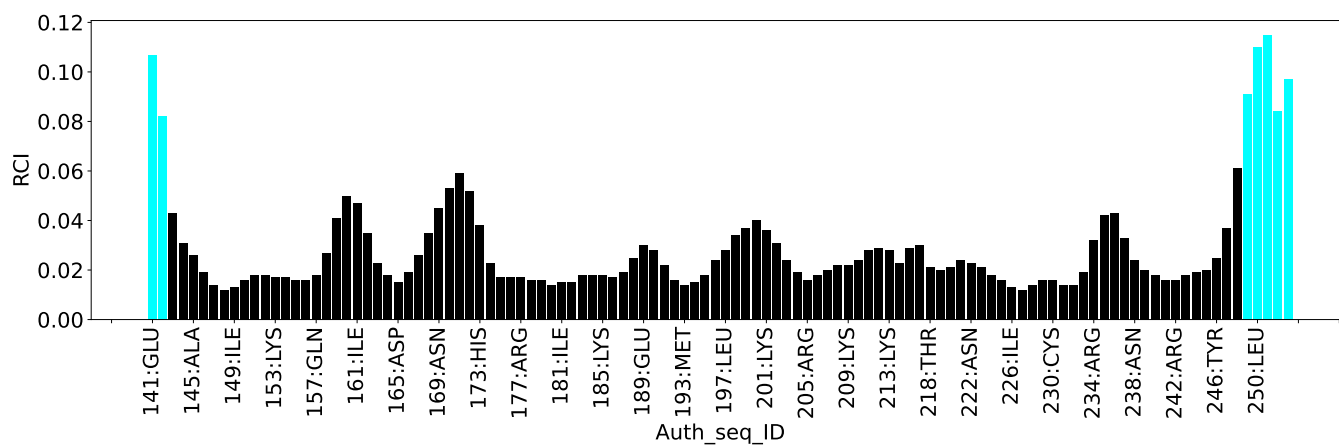
7.1.5 Random Coil Index (RCI) plots [i](#)

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

Random coil index (RCI) for chain A:



Random coil index (RCI) for chain B:



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	6432
Intra-residue ($ i-j =0$)	2067
Sequential ($ i-j =1$)	1336
Medium range ($ i-j >1$ and $ i-j <5$)	1072
Long range ($ i-j \geq 5$)	1468
Inter-chain	329
Hydrogen bond restraints	160
Disulfide bond restraints	0
Total dihedral-angle restraints	369
Number of unmapped restraints	0
Number of restraints per residue	39.8
Number of long range restraints per residue ¹	8.8

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

8.2 Residual restraint violations

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

8.2.1 Average number of distance violations per model

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

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	4.2	0.18
0.2-0.5 (Medium)	None	None
>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.

Bins (°)	Average number of violations per model	Max (°)
1.0-10.0 (Small)	4.3	4.6
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

9 Distance violation analysis [i](#)

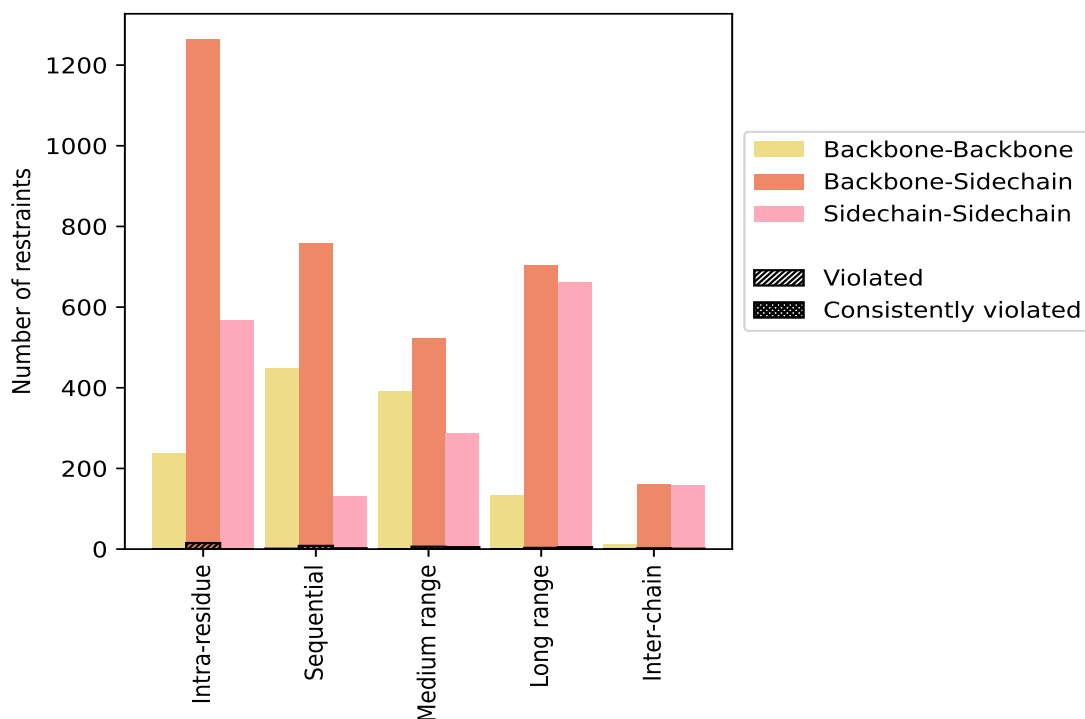
9.1 Summary of distance violations [i](#)

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

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	2067	32.1	15	0.7	0.2	0	0.0	0.0
Backbone-Backbone	237	3.7	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	1264	19.7	15	1.2	0.2	0	0.0	0.0
Sidechain-Sidechain	566	8.8	0	0.0	0.0	0	0.0	0.0
Sequential ($i-j =1$)	1336	20.8	11	0.8	0.2	0	0.0	0.0
Backbone-Backbone	447	6.9	1	0.2	0.0	0	0.0	0.0
Backbone-Sidechain	759	11.8	8	1.1	0.1	0	0.0	0.0
Sidechain-Sidechain	130	2.0	2	1.5	0.0	0	0.0	0.0
Medium range ($i-j >1$ & $i-j <5$)	1072	16.7	11	1.0	0.2	0	0.0	0.0
Backbone-Backbone	264	4.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	522	8.1	6	1.1	0.1	0	0.0	0.0
Sidechain-Sidechain	286	4.4	5	1.7	0.1	0	0.0	0.0
Long range ($i-j \geq 5$)	1468	22.8	8	0.5	0.1	0	0.0	0.0
Backbone-Backbone	102	1.6	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	704	10.9	3	0.4	0.0	0	0.0	0.0
Sidechain-Sidechain	662	10.3	5	0.8	0.1	0	0.0	0.0
Inter-chain	329	5.1	3	0.9	0.0	0	0.0	0.0
Backbone-Backbone	12	0.2	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	160	2.5	2	1.2	0.0	0	0.0	0.0
Sidechain-Sidechain	157	2.4	1	0.6	0.0	0	0.0	0.0
Hydrogen bond	160	2.5	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	6432	100.0	48	0.7	0.7	0	0.0	0.0
Backbone-Backbone	1222	19.0	1	0.1	0.0	0	0.0	0.0
Backbone-Sidechain	3409	53.0	34	1.0	0.5	0	0.0	0.0
Sidechain-Sidechain	1801	28.0	13	0.7	0.2	0	0.0	0.0

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

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



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and disulfid 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	2	1	0	0	0	3	0.15	0.15	0.0	0.15
2	2	0	0	0	0	2	0.15	0.16	0.02	0.15
3	0	2	1	1	0	4	0.13	0.14	0.01	0.13
4	0	2	0	0	0	2	0.14	0.17	0.03	0.14
5	2	0	1	0	0	3	0.13	0.16	0.02	0.12
6	1	0	0	0	1	2	0.14	0.16	0.02	0.14
7	0	1	1	1	0	3	0.13	0.14	0.01	0.14
8	1	2	1	0	0	4	0.13	0.15	0.01	0.12
9	0	1	2	0	0	3	0.13	0.16	0.02	0.12
10	3	1	0	1	0	5	0.13	0.16	0.02	0.12
11	2	3	1	1	0	7	0.12	0.16	0.02	0.12

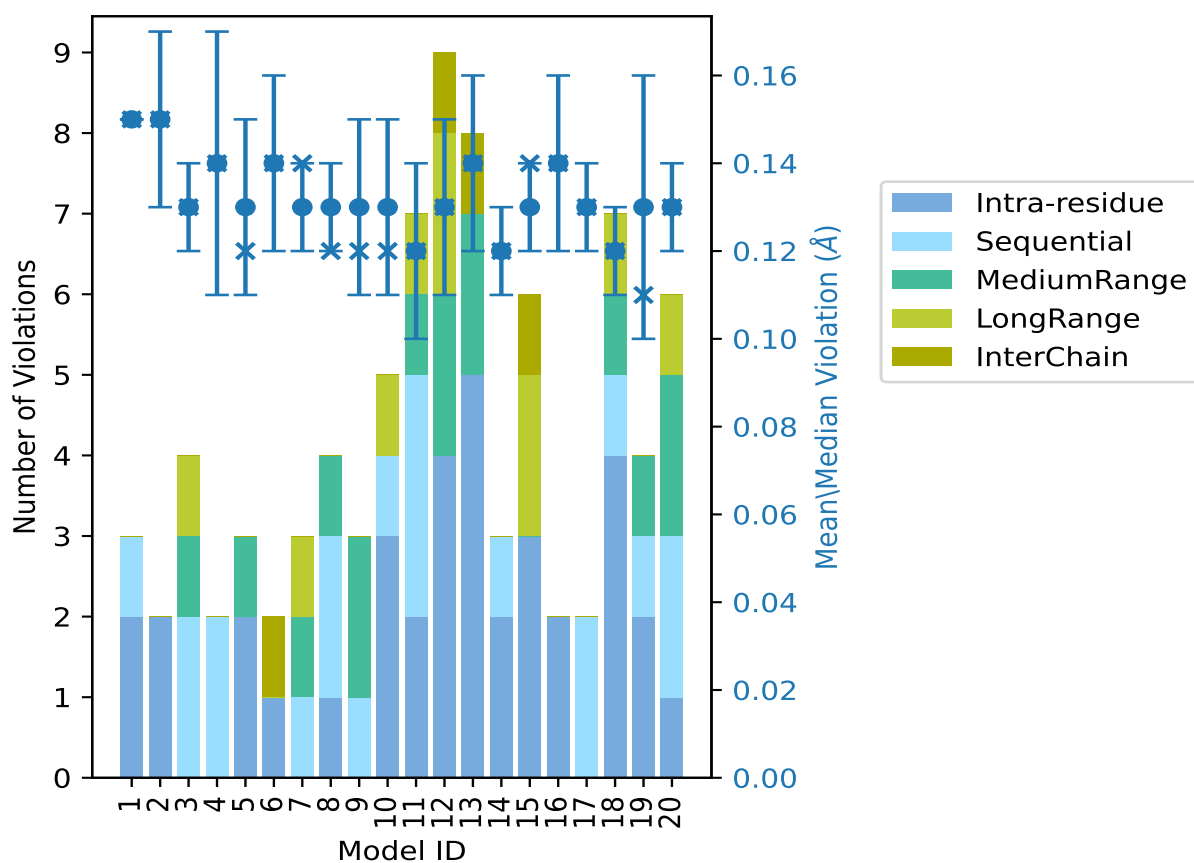
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
12	4	0	2	2	1	9	0.13	0.16	0.02	0.13
13	5	0	2	0	1	8	0.14	0.17	0.02	0.14
14	2	1	0	0	0	3	0.12	0.14	0.01	0.12
15	3	0	0	2	1	6	0.13	0.15	0.01	0.14
16	2	0	0	0	0	2	0.14	0.16	0.02	0.14
17	0	2	0	0	0	2	0.13	0.14	0.01	0.13
18	4	1	1	1	0	7	0.12	0.15	0.01	0.12
19	2	1	1	0	0	4	0.13	0.18	0.03	0.11
20	1	2	2	1	0	6	0.13	0.15	0.01	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 [\(i\)](#)



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

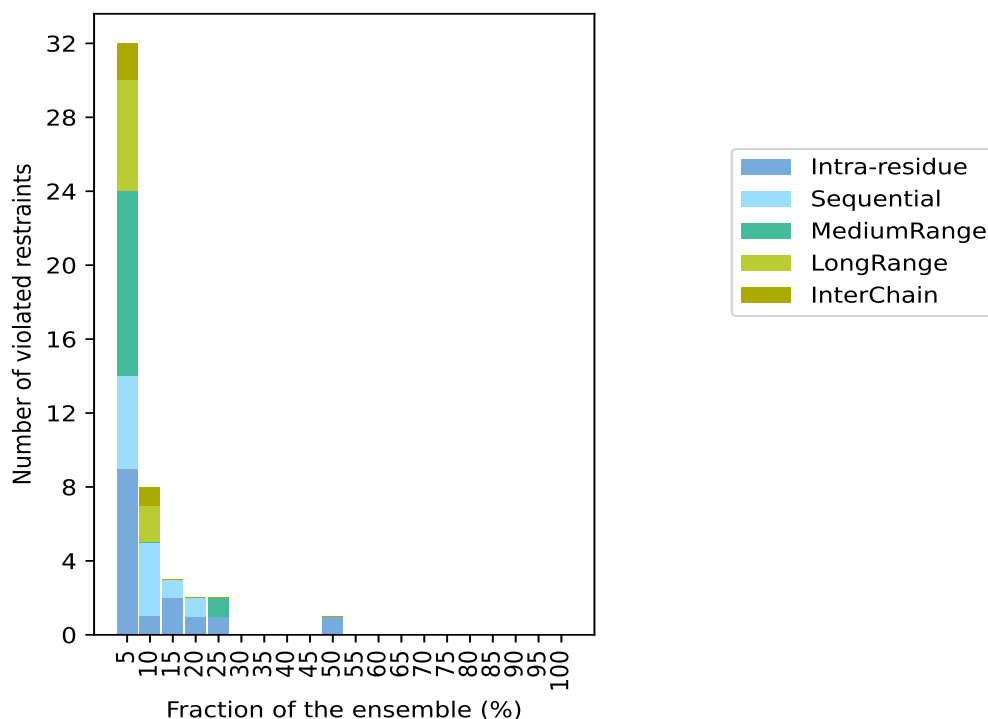
9.3 Distance violation statistics for the ensemble

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, 6224(IR:2052, SQ:1325, MR:1061, LR:1460, IC:326) 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	%
9	5	10	6	2	32	1	5.0
1	4	0	2	1	8	2	10.0
2	1	0	0	0	3	3	15.0
1	1	0	0	0	2	4	20.0
1	0	1	0	0	2	5	25.0
0	0	0	0	0	0	6	30.0
0	0	0	0	0	0	7	35.0
0	0	0	0	0	0	8	40.0
0	0	0	0	0	0	9	45.0
1	0	0	0	0	1	10	50.0
0	0	0	0	0	0	11	55.0
0	0	0	0	0	0	12	60.0
0	0	0	0	0	0	13	65.0
0	0	0	0	0	0	14	70.0
0	0	0	0	0	0	15	75.0
0	0	0	0	0	0	16	80.0
0	0	0	0	0	0	17	85.0
0	0	0	0	0	0	18	90.0
0	0	0	0	0	0	19	95.0
0	0	0	0	0	0	20	100.0

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

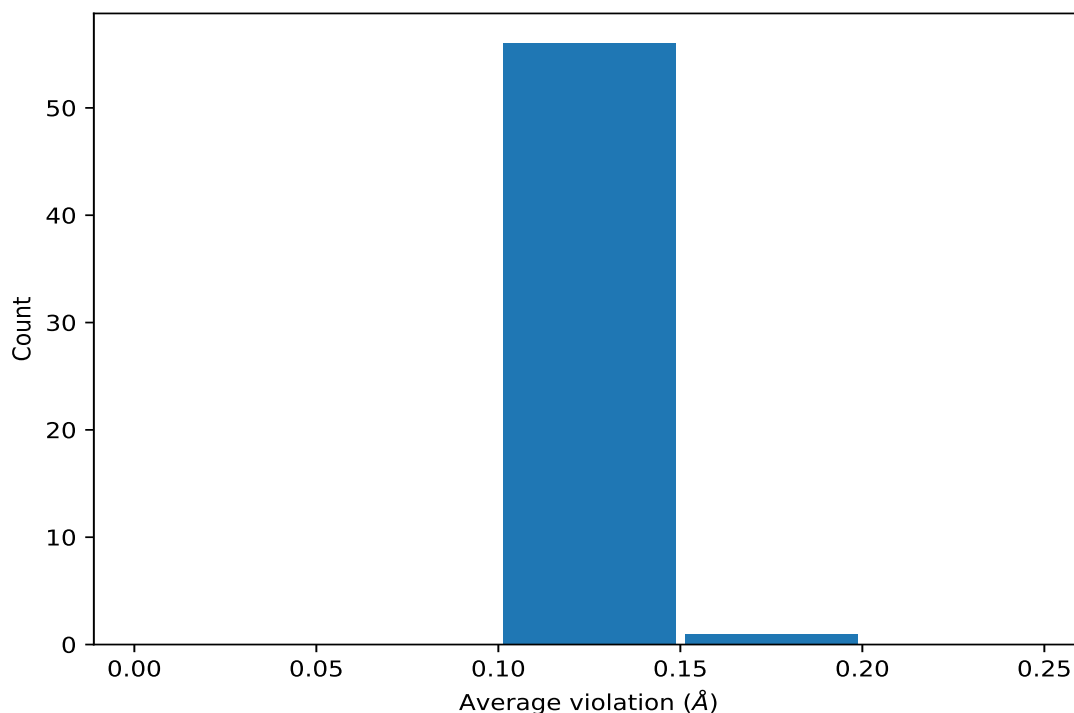
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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)
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	10	0.16	0.01	0.16
(1,507)	2:B:241:LYS:HE2	2:B:238:ASN:H	5	0.14	0.01	0.14
(1,507)	2:B:241:LYS:HE3	2:B:238:ASN:H	5	0.14	0.01	0.14
(1,3700)	2:B:251:GLU:HA	2:B:251:GLU:HG3	5	0.12	0.0	0.12
(1,4159)	1:A:50:ARG:HA	1:A:50:ARG:HB3	4	0.15	0.01	0.15
(1,1465)	1:A:44:GLU:HB2	1:A:45:ASP:H	4	0.14	0.03	0.13
(1,1465)	1:A:44:GLU:HB3	1:A:45:ASP:H	4	0.14	0.03	0.13
(1,2276)	2:B:217:ARG:HG2	2:B:218:THR:HA	3	0.12	0.01	0.12
(1,2276)	2:B:217:ARG:HG3	2:B:218:THR:HA	3	0.12	0.01	0.12
(1,3580)	2:B:151:GLU:H	2:B:151:GLU:HG2	3	0.12	0.0	0.12
(1,3580)	2:B:151:GLU:H	2:B:151:GLU:HG3	3	0.12	0.0	0.12
(1,3580)	2:B:236:ASN:HD21	2:B:151:GLU:HG2	3	0.12	0.0	0.12
(1,3580)	2:B:236:ASN:HD21	2:B:151:GLU:HG3	3	0.12	0.0	0.12
(1,3580)	2:B:154:LYS:H	2:B:151:GLU:HG2	3	0.12	0.0	0.12
(1,3580)	2:B:154:LYS:H	2:B:151:GLU:HG3	3	0.12	0.0	0.12
(1,1523)	2:B:168:LEU:HD21	2:B:168:LEU:H	3	0.12	0.0	0.12

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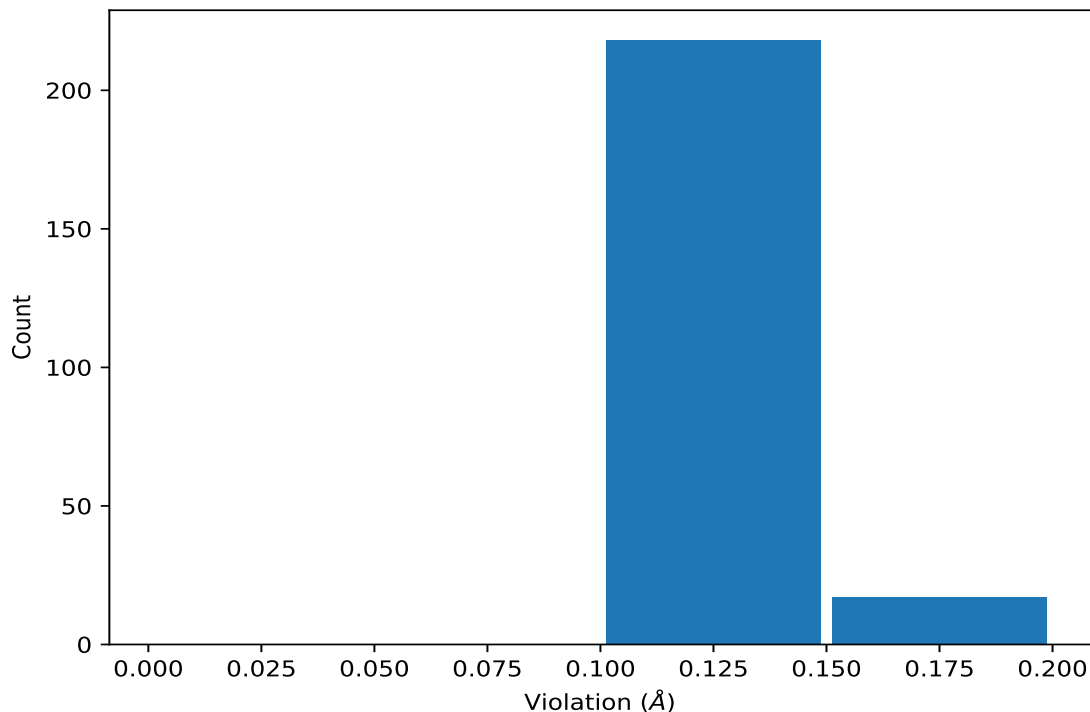
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,1523)	2:B:168:LEU:HD22	2:B:168:LEU:H	3	0.12	0.0	0.12
(1,1523)	2:B:168:LEU:HD23	2:B:168:LEU:H	3	0.12	0.0	0.12
(1,1523)	1:A:49:VAL:HG21	2:B:168:LEU:H	3	0.12	0.0	0.12
(1,1523)	1:A:49:VAL:HG22	2:B:168:LEU:H	3	0.12	0.0	0.12
(1,1523)	1:A:49:VAL:HG23	2:B:168:LEU:H	3	0.12	0.0	0.12
(1,1523)	2:B:168:LEU:HD11	2:B:168:LEU:H	3	0.12	0.0	0.12
(1,1523)	2:B:168:LEU:HD12	2:B:168:LEU:H	3	0.12	0.0	0.12
(1,1523)	2:B:168:LEU:HD13	2:B:168:LEU:H	3	0.12	0.0	0.12
(1,4189)	2:B:251:GLU:HA	2:B:251:GLU:HB3	2	0.15	0.0	0.15
(1,4189)	2:B:248:THR:HA	2:B:251:GLU:HB3	2	0.15	0.0	0.15
(1,3556)	2:B:208:LEU:H	2:B:207:GLU:HG3	2	0.14	0.01	0.14
(1,3174)	1:A:51:GLY:H	1:A:50:ARG:HD2	2	0.12	0.02	0.12
(1,3174)	1:A:51:GLY:H	1:A:50:ARG:HD3	2	0.12	0.02	0.12
(1,3278)	2:B:219:ILE:HG21	2:B:205:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HG21	2:B:205:ARG:HD3	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HG22	2:B:205:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HG22	2:B:205:ARG:HD3	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HG23	2:B:205:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HG23	2:B:205:ARG:HD3	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HD11	2:B:205:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HD11	2:B:205:ARG:HD3	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HD12	2:B:205:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HD12	2:B:205:ARG:HD3	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HD13	2:B:205:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	2:B:219:ILE:HD13	2:B:205:ARG:HD3	2	0.12	0.01	0.12
(1,3278)	1:A:35:VAL:HG11	2:B:146:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	1:A:35:VAL:HG11	2:B:146:ARG:HD3	2	0.12	0.01	0.12
(1,3278)	1:A:35:VAL:HG12	2:B:146:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	1:A:35:VAL:HG12	2:B:146:ARG:HD3	2	0.12	0.01	0.12
(1,3278)	1:A:35:VAL:HG13	2:B:146:ARG:HD2	2	0.12	0.01	0.12
(1,3278)	1:A:35:VAL:HG13	2:B:146:ARG:HD3	2	0.12	0.01	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG21	2	0.12	0.0	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG22	2	0.12	0.0	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG23	2	0.12	0.0	0.12
(1,273)	1:A:49:VAL:HG11	1:A:50:ARG:H	2	0.12	0.0	0.12
(1,273)	1:A:49:VAL:HG21	1:A:50:ARG:H	2	0.12	0.0	0.12
(1,2902)	2:B:174:ARG:HG2	2:B:175:ASP:HA	2	0.11	0.0	0.11
(1,2902)	2:B:174:ARG:HG3	2:B:175:ASP:HA	2	0.11	0.0	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG11	2	0.11	0.0	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG12	2	0.11	0.0	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG13	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)
(1,4163)	2:B:141:GLU:HA	2:B:141:GLU:HB3	19	0.18
(1,5963)	2:B:168:LEU:HG	1:A:49:VAL:HG21	13	0.17
(1,5963)	2:B:168:LEU:HG	1:A:49:VAL:HG22	13	0.17
(1,5963)	2:B:168:LEU:HG	1:A:49:VAL:HG23	13	0.17
(1,1465)	1:A:44:GLU:HB2	1:A:45:ASP:H	4	0.17
(1,1465)	1:A:44:GLU:HB3	1:A:45:ASP:H	4	0.17
(1,507)	2:B:241:LYS:HE2	2:B:238:ASN:H	9	0.16
(1,507)	2:B:241:LYS:HE3	2:B:238:ASN:H	9	0.16
(1,4159)	1:A:50:ARG:HA	1:A:50:ARG:HB3	12	0.16
(1,3928)	1:A:49:VAL:HA	1:A:49:VAL:HB	13	0.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,3928)	2:B:152:LEU:HA	2:B:155:LYS:HB3	13	0.16
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	2	0.16
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	5	0.16
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	6	0.16
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	10	0.16
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	11	0.16
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	16	0.16
(1,4189)	2:B:251:GLU:HA	2:B:251:GLU:HB3	10	0.15
(1,4189)	2:B:248:THR:HA	2:B:251:GLU:HB3	10	0.15
(1,4189)	2:B:251:GLU:HA	2:B:251:GLU:HB3	13	0.15
(1,4189)	2:B:248:THR:HA	2:B:251:GLU:HB3	13	0.15
(1,4159)	1:A:50:ARG:HA	1:A:50:ARG:HB3	1	0.15
(1,4159)	1:A:50:ARG:HA	1:A:50:ARG:HB3	18	0.15
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	12	0.15
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	15	0.15
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	20	0.15
(1,161)	1:A:52:LYS:HB2	1:A:52:LYS:H	1	0.15
(1,161)	1:A:52:LYS:HD2	1:A:52:LYS:H	1	0.15
(1,161)	1:A:52:LYS:HD3	1:A:52:LYS:H	1	0.15
(1,1465)	1:A:44:GLU:HB2	1:A:45:ASP:H	8	0.15
(1,1465)	1:A:44:GLU:HB3	1:A:45:ASP:H	8	0.15
(1,5953)	2:B:239:LEU:HD11	2:B:242:ARG:HD2	11	0.14
(1,5953)	2:B:239:LEU:HD11	2:B:242:ARG:HD3	11	0.14
(1,5953)	2:B:239:LEU:HD12	2:B:242:ARG:HD2	11	0.14
(1,5953)	2:B:239:LEU:HD12	2:B:242:ARG:HD3	11	0.14
(1,5953)	2:B:239:LEU:HD13	2:B:242:ARG:HD2	11	0.14
(1,5953)	2:B:239:LEU:HD13	2:B:242:ARG:HD3	11	0.14
(1,507)	2:B:241:LYS:HE2	2:B:238:ASN:H	3	0.14
(1,507)	2:B:241:LYS:HE3	2:B:238:ASN:H	3	0.14
(1,507)	2:B:241:LYS:HE2	2:B:238:ASN:H	7	0.14
(1,507)	2:B:241:LYS:HE3	2:B:238:ASN:H	7	0.14
(1,507)	2:B:241:LYS:HE2	2:B:238:ASN:H	12	0.14
(1,507)	2:B:241:LYS:HE3	2:B:238:ASN:H	12	0.14
(1,507)	2:B:241:LYS:HE2	2:B:238:ASN:H	13	0.14
(1,507)	2:B:241:LYS:HE3	2:B:238:ASN:H	13	0.14
(1,4560)	1:A:45:ASP:HB2	1:A:3:ARG:HG2	15	0.14
(1,4560)	1:A:45:ASP:HB2	1:A:3:ARG:HG3	15	0.14
(1,4560)	1:A:20:SER:HB3	1:A:3:ARG:HG2	15	0.14
(1,4560)	1:A:20:SER:HB3	1:A:3:ARG:HG3	15	0.14
(1,4159)	1:A:50:ARG:HA	1:A:50:ARG:HB3	13	0.14
(1,4092)	2:B:221:ASN:HD21	2:B:217:ARG:HB3	12	0.14
(1,4047)	2:B:142:ASP:H	2:B:141:GLU:HB3	7	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,3556)	2:B:208:LEU:H	2:B:207:GLU:HG3	17	0.14
(1,3174)	1:A:51:GLY:H	1:A:50:ARG:HD2	1	0.14
(1,3174)	1:A:51:GLY:H	1:A:50:ARG:HD3	1	0.14
(1,2881)	2:B:141:GLU:HB2	2:B:141:GLU:HA	14	0.14
(1,2048)	2:B:187:LEU:HD11	2:B:197:LEU:H	15	0.14
(1,2048)	2:B:187:LEU:HD12	2:B:197:LEU:H	15	0.14
(1,2048)	2:B:187:LEU:HD13	2:B:197:LEU:H	15	0.14
(1,2048)	2:B:187:LEU:HD11	2:B:208:LEU:H	15	0.14
(1,2048)	2:B:187:LEU:HD12	2:B:208:LEU:H	15	0.14
(1,2048)	2:B:187:LEU:HD13	2:B:208:LEU:H	15	0.14
(1,2048)	2:B:187:LEU:HD21	2:B:197:LEU:H	15	0.14
(1,2048)	2:B:187:LEU:HD22	2:B:197:LEU:H	15	0.14
(1,2048)	2:B:187:LEU:HD23	2:B:197:LEU:H	15	0.14
(1,2048)	2:B:214:VAL:HG21	2:B:208:LEU:H	15	0.14
(1,2048)	2:B:214:VAL:HG22	2:B:208:LEU:H	15	0.14
(1,2048)	2:B:214:VAL:HG23	2:B:208:LEU:H	15	0.14
(1,4229)	1:A:48:ASN:HB2	1:A:50:ARG:HB2	20	0.13
(1,4180)	2:B:191:GLU:HA	2:B:191:GLU:HB3	2	0.13
(1,3556)	2:B:208:LEU:H	2:B:207:GLU:HG3	20	0.13
(1,3278)	2:B:219:ILE:HG21	2:B:205:ARG:HD2	12	0.13
(1,3278)	2:B:219:ILE:HG21	2:B:205:ARG:HD3	12	0.13
(1,3278)	2:B:219:ILE:HG22	2:B:205:ARG:HD2	12	0.13
(1,3278)	2:B:219:ILE:HG22	2:B:205:ARG:HD3	12	0.13
(1,3278)	2:B:219:ILE:HG23	2:B:205:ARG:HD2	12	0.13
(1,3278)	2:B:219:ILE:HG23	2:B:205:ARG:HD3	12	0.13
(1,3278)	2:B:219:ILE:HD11	2:B:205:ARG:HD2	12	0.13
(1,3278)	2:B:219:ILE:HD11	2:B:205:ARG:HD3	12	0.13
(1,3278)	2:B:219:ILE:HD12	2:B:205:ARG:HD2	12	0.13
(1,3278)	2:B:219:ILE:HD12	2:B:205:ARG:HD3	12	0.13
(1,3278)	2:B:219:ILE:HD13	2:B:205:ARG:HD2	12	0.13
(1,3278)	2:B:219:ILE:HD13	2:B:205:ARG:HD3	12	0.13
(1,3278)	1:A:35:VAL:HG11	2:B:146:ARG:HD2	12	0.13
(1,3278)	1:A:35:VAL:HG11	2:B:146:ARG:HD3	12	0.13
(1,3278)	1:A:35:VAL:HG12	2:B:146:ARG:HD2	12	0.13
(1,3278)	1:A:35:VAL:HG12	2:B:146:ARG:HD3	12	0.13
(1,3278)	1:A:35:VAL:HG13	2:B:146:ARG:HD2	12	0.13
(1,3278)	1:A:35:VAL:HG13	2:B:146:ARG:HD3	12	0.13
(1,319)	1:A:50:ARG:H	1:A:49:VAL:H	3	0.13
(1,2949)	2:B:201:LYS:HG3	2:B:201:LYS:HA	15	0.13
(1,2949)	2:B:197:LEU:HD11	2:B:201:LYS:HA	15	0.13
(1,2949)	2:B:197:LEU:HD12	2:B:201:LYS:HA	15	0.13
(1,2949)	2:B:197:LEU:HD13	2:B:201:LYS:HA	15	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2276)	2:B:217:ARG:HG2	2:B:218:THR:HA	20	0.13
(1,2276)	2:B:217:ARG:HG3	2:B:218:THR:HA	20	0.13
(1,2150)	1:A:38:GLN:HE22	1:A:6:ILE:HD11	3	0.13
(1,2150)	1:A:38:GLN:HE22	1:A:6:ILE:HD12	3	0.13
(1,2150)	1:A:38:GLN:HE22	1:A:6:ILE:HD13	3	0.13
(1,548)	2:B:178:GLN:HB3	2:B:179:LEU:H	10	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG21	7	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG22	7	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG23	7	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG21	18	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG22	18	0.12
(1,5035)	1:A:9:GLU:HG3	1:A:35:VAL:HG23	18	0.12
(1,4539)	2:B:153:LYS:HA	2:B:231:LEU:HD21	10	0.12
(1,4539)	2:B:153:LYS:HA	2:B:231:LEU:HD22	10	0.12
(1,4539)	2:B:153:LYS:HA	2:B:231:LEU:HD23	10	0.12
(1,3713)	2:B:154:LYS:HE2	2:B:151:GLU:HG2	5	0.12
(1,3713)	2:B:154:LYS:HE2	2:B:151:GLU:HG3	5	0.12
(1,3713)	2:B:154:LYS:HE3	2:B:151:GLU:HG2	5	0.12
(1,3713)	2:B:154:LYS:HE3	2:B:151:GLU:HG3	5	0.12
(1,3713)	2:B:155:LYS:HE2	2:B:151:GLU:HG2	5	0.12
(1,3713)	2:B:155:LYS:HE2	2:B:151:GLU:HG3	5	0.12
(1,3713)	2:B:155:LYS:HE3	2:B:151:GLU:HG2	5	0.12
(1,3713)	2:B:155:LYS:HE3	2:B:151:GLU:HG3	5	0.12
(1,3700)	2:B:251:GLU:HA	2:B:251:GLU:HG3	5	0.12
(1,3700)	2:B:251:GLU:HA	2:B:251:GLU:HG3	11	0.12
(1,3700)	2:B:251:GLU:HA	2:B:251:GLU:HG3	18	0.12
(1,3580)	2:B:151:GLU:H	2:B:151:GLU:HG2	12	0.12
(1,3580)	2:B:151:GLU:H	2:B:151:GLU:HG3	12	0.12
(1,3580)	2:B:236:ASN:HD21	2:B:151:GLU:HG2	12	0.12
(1,3580)	2:B:236:ASN:HD21	2:B:151:GLU:HG3	12	0.12
(1,3580)	2:B:154:LYS:H	2:B:151:GLU:HG2	12	0.12
(1,3580)	2:B:154:LYS:H	2:B:151:GLU:HG3	12	0.12
(1,3580)	2:B:151:GLU:H	2:B:151:GLU:HG2	13	0.12
(1,3580)	2:B:151:GLU:H	2:B:151:GLU:HG3	13	0.12
(1,3580)	2:B:236:ASN:HD21	2:B:151:GLU:HG2	13	0.12
(1,3580)	2:B:236:ASN:HD21	2:B:151:GLU:HG3	13	0.12
(1,3580)	2:B:154:LYS:H	2:B:151:GLU:HG2	13	0.12
(1,3580)	2:B:154:LYS:H	2:B:151:GLU:HG3	13	0.12
(1,3580)	2:B:151:GLU:H	2:B:151:GLU:HG2	15	0.12
(1,3580)	2:B:151:GLU:H	2:B:151:GLU:HG3	15	0.12
(1,3580)	2:B:236:ASN:HD21	2:B:151:GLU:HG2	15	0.12
(1,3580)	2:B:236:ASN:HD21	2:B:151:GLU:HG3	15	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,3580)	2:B:154:LYS:H	2:B:151:GLU:HG2	15	0.12
(1,3580)	2:B:154:LYS:H	2:B:151:GLU:HG3	15	0.12
(1,3548)	1:A:50:ARG:H	1:A:48:ASN:HB2	8	0.12
(1,3278)	2:B:219:ILE:HG21	2:B:205:ARG:HD2	20	0.12
(1,3278)	2:B:219:ILE:HG21	2:B:205:ARG:HD3	20	0.12
(1,3278)	2:B:219:ILE:HG22	2:B:205:ARG:HD2	20	0.12
(1,3278)	2:B:219:ILE:HG22	2:B:205:ARG:HD3	20	0.12
(1,3278)	2:B:219:ILE:HG23	2:B:205:ARG:HD2	20	0.12
(1,3278)	2:B:219:ILE:HG23	2:B:205:ARG:HD3	20	0.12
(1,3278)	2:B:219:ILE:HD11	2:B:205:ARG:HD2	20	0.12
(1,3278)	2:B:219:ILE:HD11	2:B:205:ARG:HD3	20	0.12
(1,3278)	2:B:219:ILE:HD12	2:B:205:ARG:HD2	20	0.12
(1,3278)	2:B:219:ILE:HD12	2:B:205:ARG:HD3	20	0.12
(1,3278)	2:B:219:ILE:HD13	2:B:205:ARG:HD2	20	0.12
(1,3278)	2:B:219:ILE:HD13	2:B:205:ARG:HD3	20	0.12
(1,3278)	1:A:35:VAL:HG11	2:B:146:ARG:HD2	20	0.12
(1,3278)	1:A:35:VAL:HG11	2:B:146:ARG:HD3	20	0.12
(1,3278)	1:A:35:VAL:HG12	2:B:146:ARG:HD2	20	0.12
(1,3278)	1:A:35:VAL:HG12	2:B:146:ARG:HD3	20	0.12
(1,3278)	1:A:35:VAL:HG13	2:B:146:ARG:HD2	20	0.12
(1,3278)	1:A:35:VAL:HG13	2:B:146:ARG:HD3	20	0.12
(1,3249)	2:B:158:GLU:HB2	2:B:201:LYS:HE2	12	0.12
(1,3195)	2:B:221:ASN:HD22	2:B:217:ARG:HD2	9	0.12
(1,2903)	2:B:168:LEU:HG	2:B:165:ASP:HA	13	0.12
(1,2884)	1:A:1:MET:HG2	1:A:1:MET:HA	8	0.12
(1,2884)	1:A:1:MET:HG3	1:A:1:MET:HA	8	0.12
(1,273)	1:A:49:VAL:HG11	1:A:50:ARG:H	17	0.12
(1,273)	1:A:49:VAL:HG21	1:A:50:ARG:H	17	0.12
(1,2658)	2:B:155:LYS:HG2	2:B:155:LYS:HA	18	0.12
(1,2658)	2:B:154:LYS:HG3	2:B:155:LYS:HA	18	0.12
(1,2276)	2:B:217:ARG:HG2	2:B:218:THR:HA	8	0.12
(1,2276)	2:B:217:ARG:HG3	2:B:218:THR:HA	8	0.12
(1,1906)	2:B:209:LYS:HG2	2:B:214:VAL:H	11	0.12
(1,1906)	2:B:209:LYS:HG3	2:B:214:VAL:H	11	0.12
(1,1906)	2:B:219:ILE:HG13	2:B:214:VAL:H	11	0.12
(1,1523)	2:B:168:LEU:HD21	2:B:168:LEU:H	13	0.12
(1,1523)	2:B:168:LEU:HD22	2:B:168:LEU:H	13	0.12
(1,1523)	2:B:168:LEU:HD23	2:B:168:LEU:H	13	0.12
(1,1523)	1:A:49:VAL:HG21	2:B:168:LEU:H	13	0.12
(1,1523)	1:A:49:VAL:HG22	2:B:168:LEU:H	13	0.12
(1,1523)	1:A:49:VAL:HG23	2:B:168:LEU:H	13	0.12
(1,1523)	2:B:168:LEU:HD11	2:B:168:LEU:H	13	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1523)	2:B:168:LEU:HD12	2:B:168:LEU:H	13	0.12
(1,1523)	2:B:168:LEU:HD13	2:B:168:LEU:H	13	0.12
(1,1523)	2:B:168:LEU:HD21	2:B:168:LEU:H	14	0.12
(1,1523)	2:B:168:LEU:HD22	2:B:168:LEU:H	14	0.12
(1,1523)	2:B:168:LEU:HD23	2:B:168:LEU:H	14	0.12
(1,1523)	1:A:49:VAL:HG21	2:B:168:LEU:H	14	0.12
(1,1523)	1:A:49:VAL:HG22	2:B:168:LEU:H	14	0.12
(1,1523)	1:A:49:VAL:HG23	2:B:168:LEU:H	14	0.12
(1,1523)	2:B:168:LEU:HD11	2:B:168:LEU:H	14	0.12
(1,1523)	2:B:168:LEU:HD12	2:B:168:LEU:H	14	0.12
(1,1523)	2:B:168:LEU:HD13	2:B:168:LEU:H	14	0.12
(1,952)	2:B:153:LYS:HE2	2:B:153:LYS:H	16	0.11
(1,952)	2:B:153:LYS:HE3	2:B:153:LYS:H	16	0.11
(1,952)	2:B:154:LYS:HE2	2:B:153:LYS:H	16	0.11
(1,952)	2:B:154:LYS:HE3	2:B:153:LYS:H	16	0.11
(1,5967)	2:B:172:LYS:HE2	2:B:173:HIS:HE1	11	0.11
(1,5967)	2:B:172:LYS:HE3	2:B:173:HIS:HE1	11	0.11
(1,5950)	2:B:241:LYS:HG3	2:B:238:ASN:HA	19	0.11
(1,593)	2:B:154:LYS:HG2	2:B:151:GLU:H	18	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG11	12	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG12	12	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG13	12	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG11	15	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG12	15	0.11
(1,5915)	2:B:168:LEU:HA	1:A:49:VAL:HG13	15	0.11
(1,5737)	1:A:23:GLU:HG2	2:B:171:PRO:HA	6	0.11
(1,3743)	2:B:149:ILE:HG12	2:B:150:GLU:HG2	19	0.11
(1,3743)	2:B:149:ILE:HG12	2:B:147:GLU:HG2	19	0.11
(1,3743)	2:B:154:LYS:HG2	2:B:150:GLU:HG2	19	0.11
(1,3700)	2:B:251:GLU:HA	2:B:251:GLU:HG3	12	0.11
(1,3700)	2:B:251:GLU:HA	2:B:251:GLU:HG3	19	0.11
(1,3174)	1:A:51:GLY:H	1:A:50:ARG:HD2	14	0.11
(1,3174)	1:A:51:GLY:H	1:A:50:ARG:HD3	14	0.11
(1,2902)	2:B:174:ARG:HG2	2:B:175:ASP:HA	11	0.11
(1,2902)	2:B:174:ARG:HG3	2:B:175:ASP:HA	11	0.11
(1,2902)	2:B:174:ARG:HG2	2:B:175:ASP:HA	18	0.11
(1,2902)	2:B:174:ARG:HG3	2:B:175:ASP:HA	18	0.11
(1,273)	1:A:49:VAL:HG11	1:A:50:ARG:H	9	0.11
(1,273)	1:A:49:VAL:HG21	1:A:50:ARG:H	9	0.11
(1,2619)	2:B:209:LYS:HE2	2:B:209:LYS:HA	10	0.11
(1,2619)	2:B:209:LYS:HE3	2:B:209:LYS:HA	10	0.11
(1,2276)	2:B:217:ARG:HG2	2:B:218:THR:HA	4	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,2276)	2:B:217:ARG:HG3	2:B:218:THR:HA	4	0.11
(1,1523)	2:B:168:LEU:HD21	2:B:168:LEU:H	18	0.11
(1,1523)	2:B:168:LEU:HD22	2:B:168:LEU:H	18	0.11
(1,1523)	2:B:168:LEU:HD23	2:B:168:LEU:H	18	0.11
(1,1523)	1:A:49:VAL:HG21	2:B:168:LEU:H	18	0.11
(1,1523)	1:A:49:VAL:HG22	2:B:168:LEU:H	18	0.11
(1,1523)	1:A:49:VAL:HG23	2:B:168:LEU:H	18	0.11
(1,1523)	2:B:168:LEU:HD11	2:B:168:LEU:H	18	0.11
(1,1523)	2:B:168:LEU:HD12	2:B:168:LEU:H	18	0.11
(1,1523)	2:B:168:LEU:HD13	2:B:168:LEU:H	18	0.11
(1,1465)	1:A:44:GLU:HB2	1:A:45:ASP:H	3	0.11
(1,1465)	1:A:44:GLU:HB3	1:A:45:ASP:H	3	0.11
(1,1465)	1:A:44:GLU:HB2	1:A:45:ASP:H	11	0.11
(1,1465)	1:A:44:GLU:HB3	1:A:45:ASP:H	11	0.11
(1,1018)	2:B:207:GLU:HG2	2:B:211:LYS:H	20	0.11

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

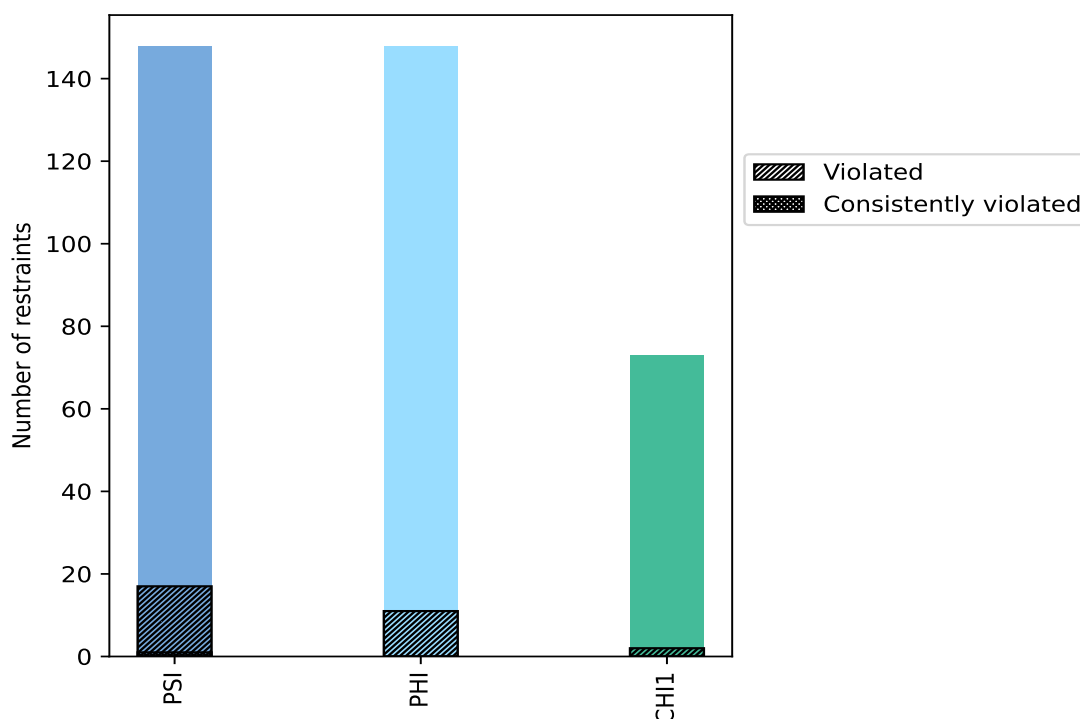
10.1 Summary of dihedral-angle violations [i](#)

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

Angle type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
PSI	148	40.1	17	11.5	4.6	1	0.7	0.3
PHI	148	40.1	11	7.4	3.0	0	0.0	0.0
CHI1	73	19.8	2	2.7	0.5	0	0.0	0.0
Total	369	100.0	30	8.1	8.1	1	0.3	0.3

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

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



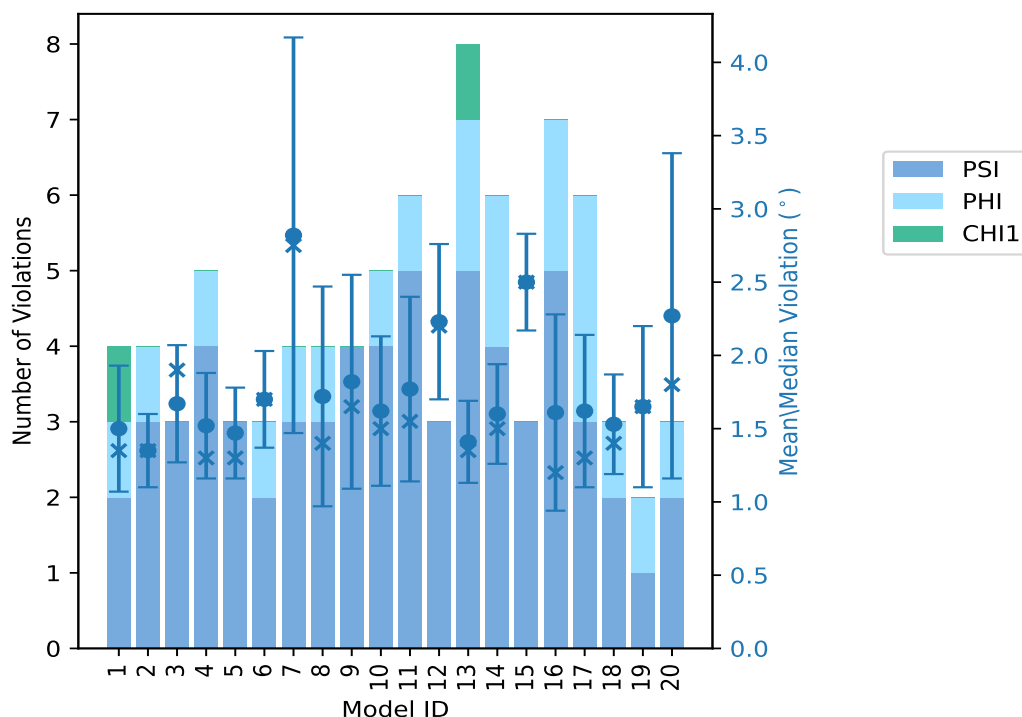
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

10.2 Dihedral-angle violation statistics for each model [i](#)

The following table provides the dihedral-angle violation statistics for each model in the ensemble. Violations less than 1° are not included in the statistics.

Model ID	Number of violations				Mean (°)	Max (°)	SD (°)	Median (°)
	PSI	PHI	CHI1	Total				
1	2	1	1	4	1.5	2.2	0.43	1.35
2	3	1	0	4	1.35	1.6	0.25	1.35
3	3	0	0	3	1.67	2.0	0.4	1.9
4	4	1	0	5	1.52	2.0	0.36	1.3
5	3	0	0	3	1.47	1.9	0.31	1.3
6	2	1	0	3	1.7	2.1	0.33	1.7
7	3	1	0	4	2.82	4.6	1.35	2.75
8	3	1	0	4	1.72	3.0	0.75	1.4
9	4	0	0	4	1.82	2.9	0.73	1.65
10	4	1	0	5	1.62	2.4	0.51	1.5
11	5	1	0	6	1.77	3.1	0.63	1.55
12	3	0	0	3	2.23	2.9	0.53	2.2
13	5	2	1	8	1.41	2.0	0.28	1.35
14	4	2	0	6	1.6	2.1	0.34	1.5
15	3	0	0	3	2.5	2.9	0.33	2.5
16	5	2	0	7	1.61	3.0	0.67	1.2
17	3	3	0	6	1.62	2.4	0.52	1.3
18	2	1	0	3	1.53	2.0	0.34	1.4
19	1	1	0	2	1.65	2.2	0.55	1.65
20	2	1	0	3	2.27	3.8	1.11	1.8

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



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

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

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

Number of violated restraints				Fraction of the ensemble	
PSI	PHI	CHI1	Total	Count ¹	%
8	5	2	15	1	5.0
3	3	0	6	2	10.0
4	3	0	7	3	15.0
0	0	0	0	4	20.0
0	0	0	0	5	25.0
0	0	0	0	6	30.0
0	0	0	0	7	35.0
0	0	0	0	8	40.0
0	0	0	0	9	45.0
0	0	0	0	10	50.0
0	0	0	0	11	55.0

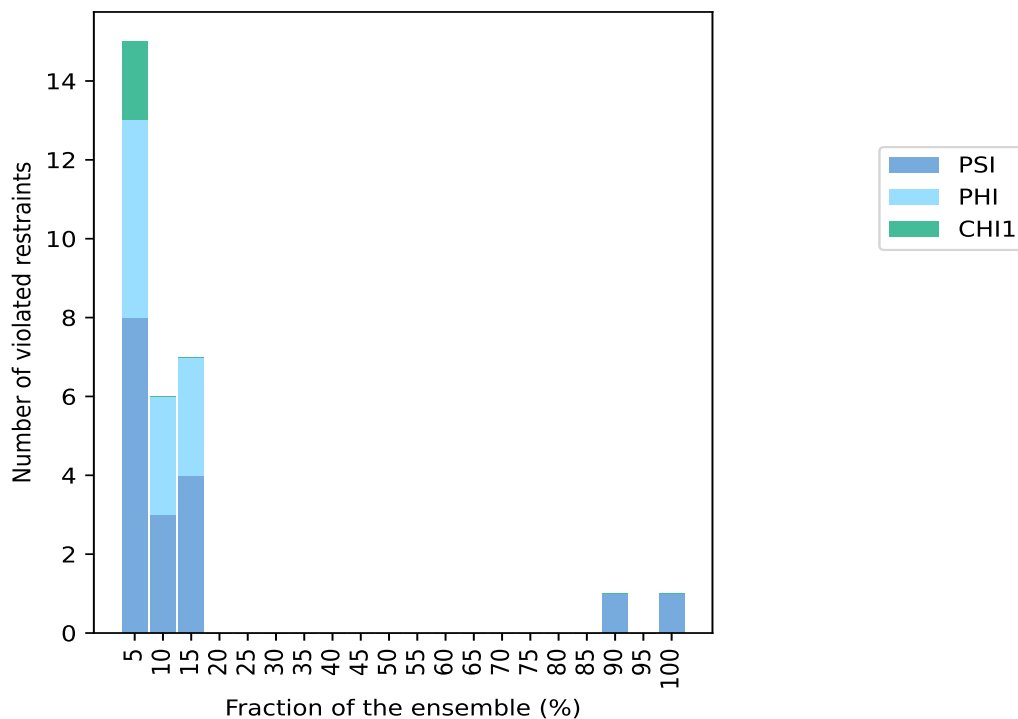
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Number of violated restraints				Fraction of the ensemble	
PSI	PHI	CHI1	Total	Count ¹	%
0	0	0	0	12	60.0
0	0	0	0	13	65.0
0	0	0	0	14	70.0
0	0	0	0	15	75.0
0	0	0	0	16	80.0
0	0	0	0	17	85.0
1	0	0	1	18	90.0
0	0	0	0	19	95.0
1	0	0	1	20	100.0

¹ Number of models with violations

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

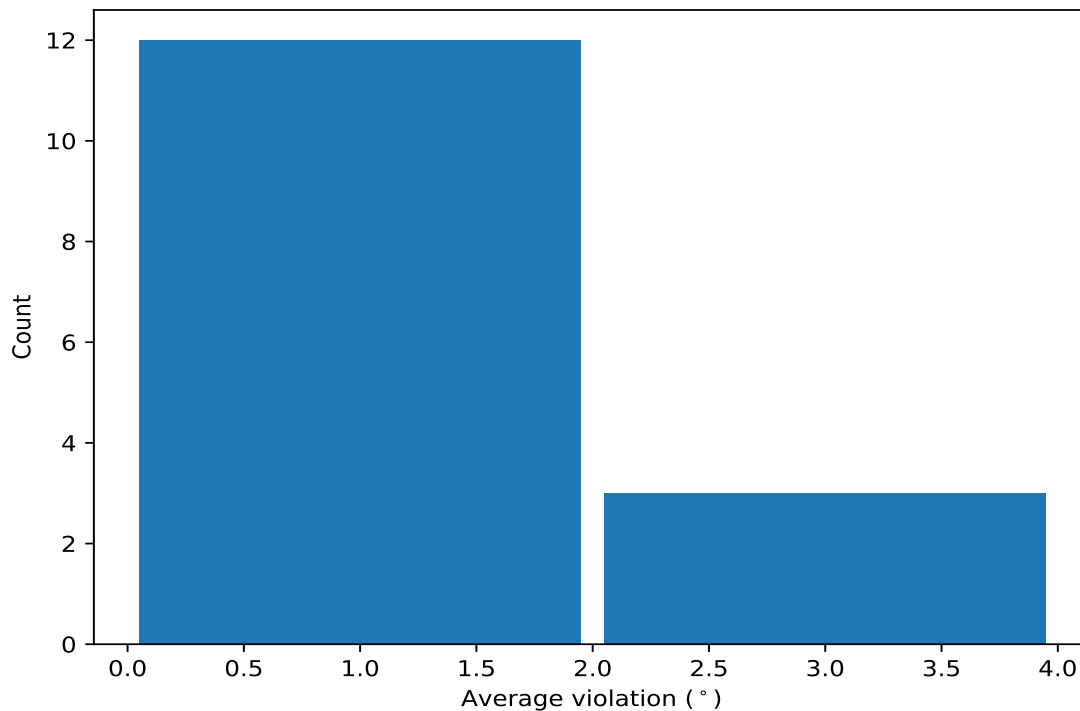


10.4 Most violated dihedral-angle restraints in the ensemble [i](#)

10.4.1 Histogram : Distribution of mean dihedral-angle violations [i](#)

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

in the ensemble



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

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

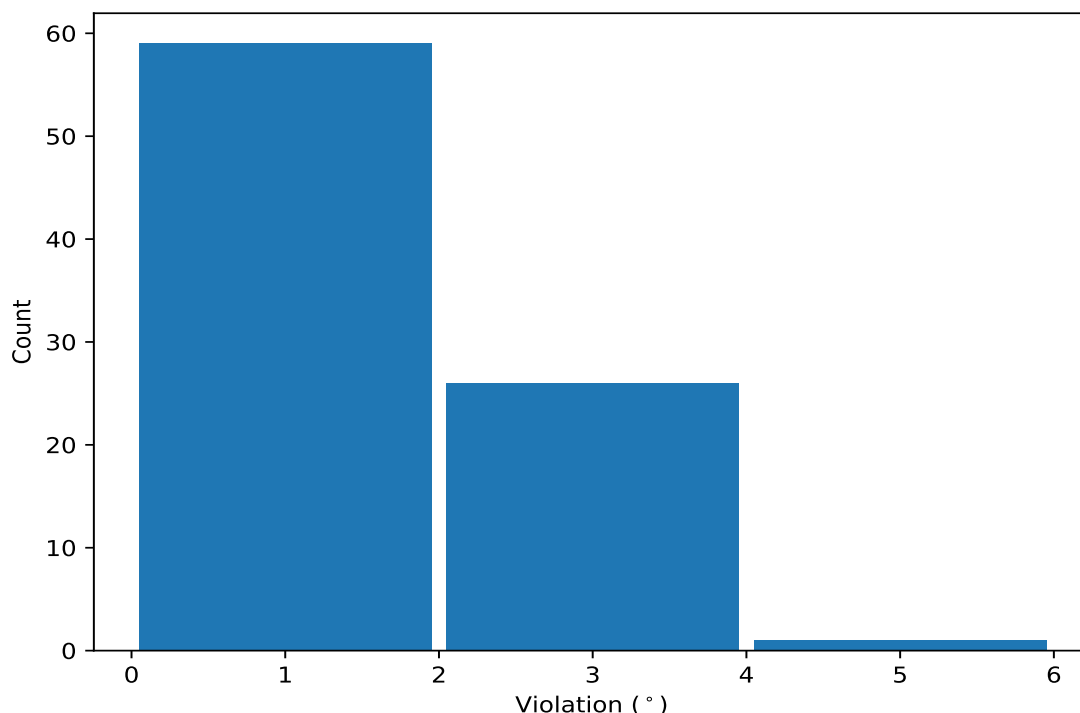
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models ¹	Mean	SD ²	Median
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	20	2.26	0.71	2.05
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	18	2.11	0.43	2.05
(1,58)	1:A:32:ASP:N	1:A:32:ASP:CA	1:A:32:ASP:C	1:A:33:MET:N	3	1.47	0.45	1.2
(1,61)	1:A:33:MET:C	1:A:34:LYS:N	1:A:34:LYS:CA	1:A:34:LYS:C	3	1.43	0.09	1.5
(1,231)	2:B:214:VAL:C	2:B:215:HIS:N	2:B:215:HIS:CA	2:B:215:HIS:C	3	1.4	0.28	1.2
(1,1)	1:A:1:MET:C	1:A:2:ASN:N	1:A:2:ASN:CA	1:A:2:ASN:C	3	1.23	0.19	1.1
(1,180)	2:B:189:GLU:N	2:B:189:GLU:CA	2:B:189:GLU:C	2:B:190:ASP:N	3	1.23	0.19	1.1
(1,32)	1:A:18:LEU:N	1:A:18:LEU:CA	1:A:18:LEU:C	1:A:19:THR:N	3	1.2	0.08	1.2
(1,18)	1:A:11:GLN:N	1:A:11:GLN:CA	1:A:11:GLN:C	1:A:12:GLY:N	3	1.17	0.05	1.2
(1,296)	2:B:256:HIS:N	2:B:256:HIS:CA	2:B:256:HIS:C	2:B:257:HIS:N	2	3.1	1.5	3.1
(1,239)	2:B:219:ILE:C	2:B:220:GLY:N	2:B:220:GLY:CA	2:B:220:GLY:C	2	1.45	0.05	1.45
(1,91)	2:B:144:GLU:C	2:B:145:ALA:N	2:B:145:ALA:CA	2:B:145:ALA:C	2	1.3	0.0	1.3
(1,102)	2:B:150:GLU:N	2:B:150:GLU:CA	2:B:150:GLU:C	2:B:151:GLU:N	2	1.25	0.15	1.25
(1,174)	2:B:186:MET:N	2:B:186:MET:CA	2:B:186:MET:C	2:B:187:LEU:N	2	1.25	0.05	1.25
(1,75)	1:A:40:SER:C	1:A:41:PHE:N	1:A:41:PHE:CA	1:A:41:PHE:C	2	1.1	0.0	1.1

¹ Number of violated models, ²Standard deviation, All angle values are in degree (°)

10.5 All violated dihedral-angle restraints [i](#)

10.5.1 Histogram : Distribution of violations [i](#)

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



10.5.2 Table: All violated dihedral-angle restraints [i](#)

The following table lists the absolute value of the violation for each restraint in the ensemble sorted by its value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint.

Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,296)	2:B:256:HIS:N	2:B:256:HIS:CA	2:B:256:HIS:C	2:B:257:HIS:N	7	4.6
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	20	3.8
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	7	3.6
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	11	3.1
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	8	3.0
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	16	3.0
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	9	2.9
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	15	2.9
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	12	2.9
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	15	2.5
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	10	2.4
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	17	2.4
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	17	2.3
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	1	2.2

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	12	2.2
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	16	2.2
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	19	2.2
(1,58)	1:A:32:ASP:N	1:A:32:ASP:CA	1:A:32:ASP:C	1:A:33:MET:N	15	2.1
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	6	2.1
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	14	2.1
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	9	2.1
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	3	2.0
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	10	2.0
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	4	2.0
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	13	2.0
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	14	2.0
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	18	2.0
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	4	1.9
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	5	1.9
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	7	1.9
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	3	1.9
(1,248)	2:B:224:LYS:N	2:B:224:LYS:CA	2:B:224:LYS:C	2:B:225:TYR:N	11	1.8
(1,231)	2:B:214:VAL:C	2:B:215:HIS:N	2:B:215:HIS:CA	2:B:215:HIS:C	20	1.8
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	11	1.7
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	13	1.7
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	6	1.7
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	2	1.6
(1,296)	2:B:256:HIS:N	2:B:256:HIS:CA	2:B:256:HIS:C	2:B:257:HIS:N	12	1.6
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	2	1.6
(1,61)	1:A:33:MET:C	1:A:34:LYS:N	1:A:34:LYS:CA	1:A:34:LYS:C	10	1.5
(1,61)	1:A:33:MET:C	1:A:34:LYS:N	1:A:34:LYS:CA	1:A:34:LYS:C	14	1.5
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	1	1.5
(1,239)	2:B:219:ILE:C	2:B:220:GLY:N	2:B:220:GLY:CA	2:B:220:GLY:C	14	1.5
(1,180)	2:B:189:GLU:N	2:B:189:GLU:CA	2:B:189:GLU:C	2:B:190:ASP:N	8	1.5
(1,1)	1:A:1:MET:C	1:A:2:ASN:N	1:A:2:ASN:CA	1:A:2:ASN:C	16	1.5
(1,42)	1:A:24:PHE:N	1:A:24:PHE:CA	1:A:24:PHE:C	1:A:25:LEU:N	18	1.4
(1,239)	2:B:219:ILE:C	2:B:220:GLY:N	2:B:220:GLY:CA	2:B:220:GLY:C	17	1.4
(1,150)	2:B:174:ARG:N	2:B:174:ARG:CA	2:B:174:ARG:C	2:B:175:ASP:N	13	1.4
(1,149)	2:B:173:HIS:C	2:B:174:ARG:N	2:B:174:ARG:CA	2:B:174:ARG:C	13	1.4
(1,102)	2:B:150:GLU:N	2:B:150:GLU:CA	2:B:150:GLU:C	2:B:151:GLU:N	11	1.4
(1,91)	2:B:144:GLU:C	2:B:145:ALA:N	2:B:145:ALA:CA	2:B:145:ALA:C	8	1.3
(1,91)	2:B:144:GLU:C	2:B:145:ALA:N	2:B:145:ALA:CA	2:B:145:ALA:C	11	1.3
(1,74)	1:A:40:SER:N	1:A:40:SER:CA	1:A:40:SER:C	1:A:41:PHE:N	13	1.3
(1,61)	1:A:33:MET:C	1:A:34:LYS:N	1:A:34:LYS:CA	1:A:34:LYS:C	6	1.3
(1,32)	1:A:18:LEU:N	1:A:18:LEU:CA	1:A:18:LEU:C	1:A:19:THR:N	14	1.3
(1,271)	2:B:237:LEU:C	2:B:238:ASN:N	2:B:238:ASN:CA	2:B:238:ASN:C	4	1.3
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	5	1.3
(1,242)	2:B:221:ASN:N	2:B:221:ASN:CA	2:B:221:ASN:C	2:B:222:ASN:N	13	1.3
(1,24)	1:A:14:LYS:N	1:A:14:LYS:CA	1:A:14:LYS:C	1:A:15:ALA:N	4	1.3
(1,174)	2:B:186:MET:N	2:B:186:MET:CA	2:B:186:MET:C	2:B:187:LEU:N	11	1.3
(1,70)	1:A:38:GLN:N	1:A:38:GLN:CA	1:A:38:GLN:C	1:A:39:VAL:N	16	1.2
(1,58)	1:A:32:ASP:N	1:A:32:ASP:CA	1:A:32:ASP:C	1:A:33:MET:N	14	1.2
(1,349)	2:B:202:ASN:N	2:B:202:ASN:CA	2:B:202:ASN:CB	2:B:202:ASN:CG	1	1.2
(1,32)	1:A:18:LEU:N	1:A:18:LEU:CA	1:A:18:LEU:C	1:A:19:THR:N	16	1.2
(1,295)	2:B:255:HIS:C	2:B:256:HIS:N	2:B:256:HIS:CA	2:B:256:HIS:C	17	1.2

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Key	Atom-1	Atom-2	Atom-3	Atom-4	Model ID	Violation (°)
(1,231)	2:B:214:VAL:C	2:B:215:HIS:N	2:B:215:HIS:CA	2:B:215:HIS:C	7	1.2
(1,231)	2:B:214:VAL:C	2:B:215:HIS:N	2:B:215:HIS:CA	2:B:215:HIS:C	18	1.2
(1,18)	1:A:11:GLN:N	1:A:11:GLN:CA	1:A:11:GLN:C	1:A:12:GLY:N	5	1.2
(1,18)	1:A:11:GLN:N	1:A:11:GLN:CA	1:A:11:GLN:C	1:A:12:GLY:N	17	1.2
(1,179)	2:B:188:ALA:C	2:B:189:GLU:N	2:B:189:GLU:CA	2:B:189:GLU:C	17	1.2
(1,174)	2:B:186:MET:N	2:B:186:MET:CA	2:B:186:MET:C	2:B:187:LEU:N	9	1.2
(1,120)	2:B:159:PHE:N	2:B:159:PHE:CA	2:B:159:PHE:C	2:B:160:GLY:N	20	1.2
(1,75)	1:A:40:SER:C	1:A:41:PHE:N	1:A:41:PHE:CA	1:A:41:PHE:C	2	1.1
(1,75)	1:A:40:SER:C	1:A:41:PHE:N	1:A:41:PHE:CA	1:A:41:PHE:C	16	1.1
(1,58)	1:A:32:ASP:N	1:A:32:ASP:CA	1:A:32:ASP:C	1:A:33:MET:N	10	1.1
(1,357)	2:B:221:ASN:N	2:B:221:ASN:CA	2:B:221:ASN:CB	2:B:221:ASN:CG	13	1.1
(1,32)	1:A:18:LEU:N	1:A:18:LEU:CA	1:A:18:LEU:C	1:A:19:THR:N	2	1.1
(1,270)	2:B:235:SER:N	2:B:235:SER:CA	2:B:235:SER:C	2:B:236:ASN:N	8	1.1
(1,230)	2:B:214:VAL:N	2:B:214:VAL:CA	2:B:214:VAL:C	2:B:215:HIS:N	9	1.1
(1,229)	2:B:213:LYS:C	2:B:214:VAL:N	2:B:214:VAL:CA	2:B:214:VAL:C	13	1.1
(1,180)	2:B:189:GLU:N	2:B:189:GLU:CA	2:B:189:GLU:C	2:B:190:ASP:N	3	1.1
(1,180)	2:B:189:GLU:N	2:B:189:GLU:CA	2:B:189:GLU:C	2:B:190:ASP:N	16	1.1
(1,18)	1:A:11:GLN:N	1:A:11:GLN:CA	1:A:11:GLN:C	1:A:12:GLY:N	4	1.1
(1,102)	2:B:150:GLU:N	2:B:150:GLU:CA	2:B:150:GLU:C	2:B:151:GLU:N	10	1.1
(1,1)	1:A:1:MET:C	1:A:2:ASN:N	1:A:2:ASN:CA	1:A:2:ASN:C	1	1.1
(1,1)	1:A:1:MET:C	1:A:2:ASN:N	1:A:2:ASN:CA	1:A:2:ASN:C	19	1.1