



Full wwPDB NMR Structure Validation Report ⓘ

Jun 3, 2023 – 05:54 PM EDT

PDB ID : 2MP1
BMRB ID : 19960
Title : Solution structure of the human chemokine CCL19
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Deposited on : 2014-05-09

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

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with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-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

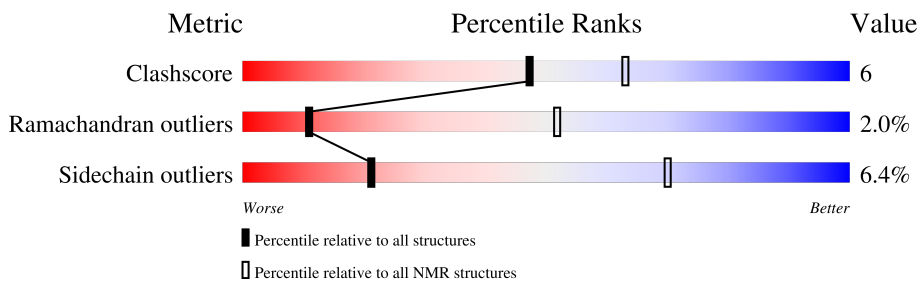
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 83%.

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	77	

2 Ensemble composition and analysis

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

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

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:8-A:67 (60)	0.63	1

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

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

The models can be grouped into 4 clusters and 3 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1250 atoms, of which 634 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called C-C motif chemokine 19.

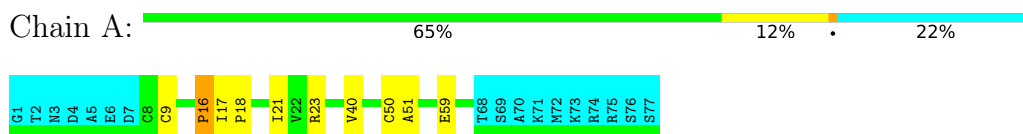
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	77	1250	385	634	118	108	5	0

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: C-C motif chemokine 19

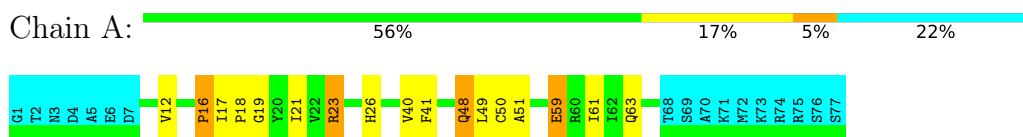


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

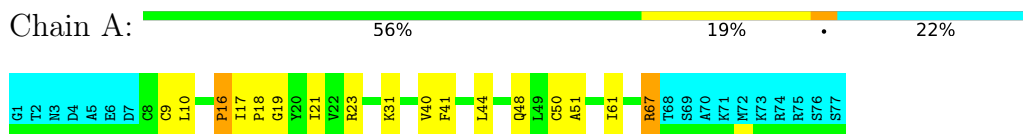
4.2.1 Score per residue for model 1 (medoid)

- Molecule 1: C-C motif chemokine 19



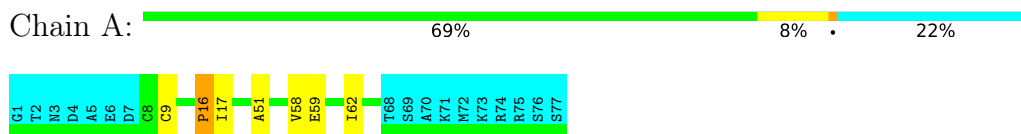
4.2.2 Score per residue for model 2

- Molecule 1: C-C motif chemokine 19



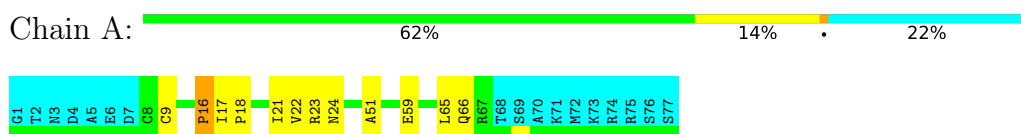
4.2.3 Score per residue for model 3

- Molecule 1: C-C motif chemokine 19



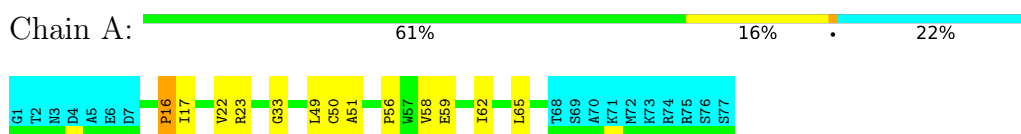
4.2.4 Score per residue for model 4

- Molecule 1: C-C motif chemokine 19



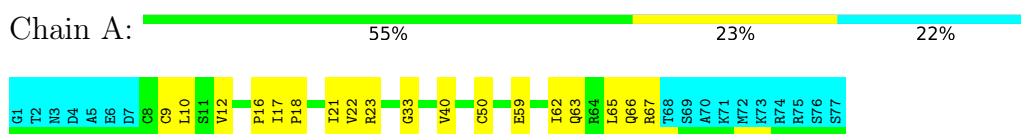
4.2.5 Score per residue for model 5

- Molecule 1: C-C motif chemokine 19



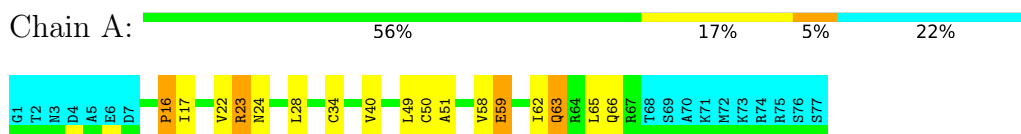
4.2.6 Score per residue for model 6

- Molecule 1: C-C motif chemokine 19



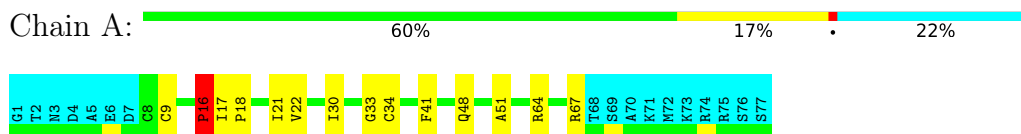
4.2.7 Score per residue for model 7

- Molecule 1: C-C motif chemokine 19



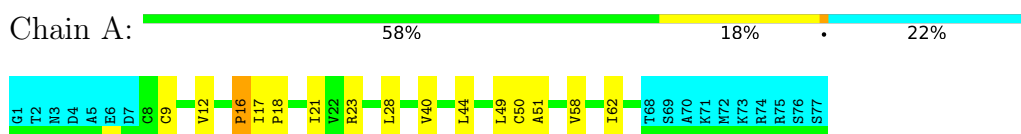
4.2.8 Score per residue for model 8

- Molecule 1: C-C motif chemokine 19



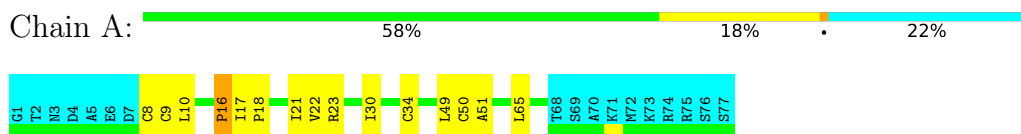
4.2.9 Score per residue for model 9

- Molecule 1: C-C motif chemokine 19



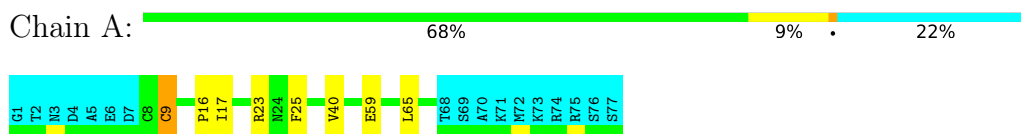
4.2.10 Score per residue for model 10

- Molecule 1: C-C motif chemokine 19



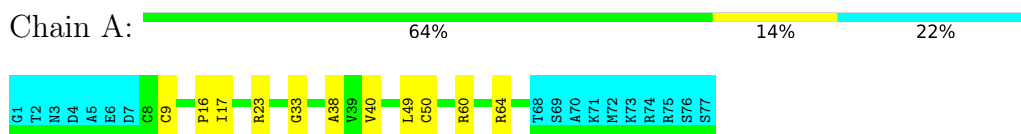
4.2.11 Score per residue for model 11

- Molecule 1: C-C motif chemokine 19



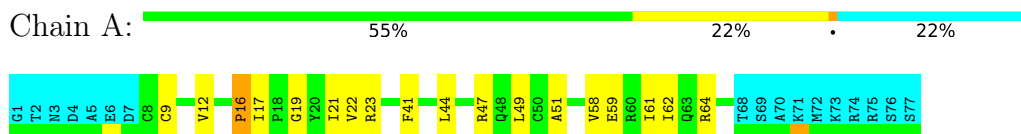
4.2.12 Score per residue for model 12

- Molecule 1: C-C motif chemokine 19



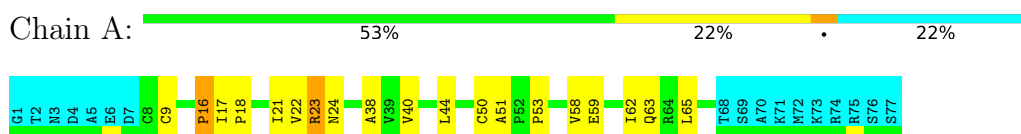
4.2.13 Score per residue for model 13

- Molecule 1: C-C motif chemokine 19



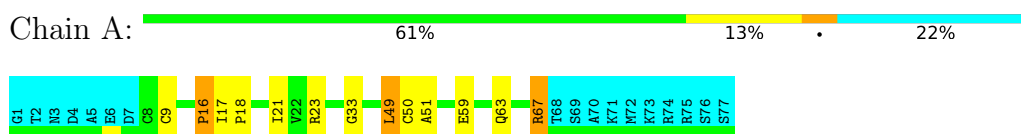
4.2.14 Score per residue for model 14

- Molecule 1: C-C motif chemokine 19



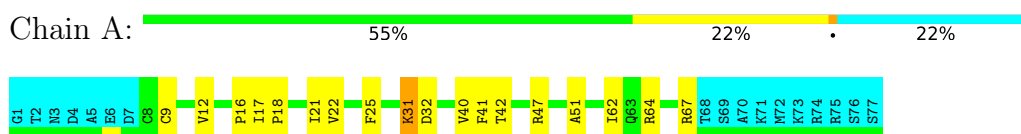
4.2.15 Score per residue for model 15

- Molecule 1: C-C motif chemokine 19



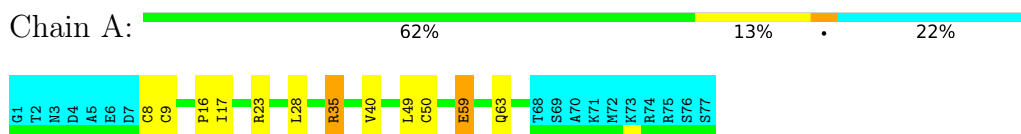
4.2.16 Score per residue for model 16

- Molecule 1: C-C motif chemokine 19



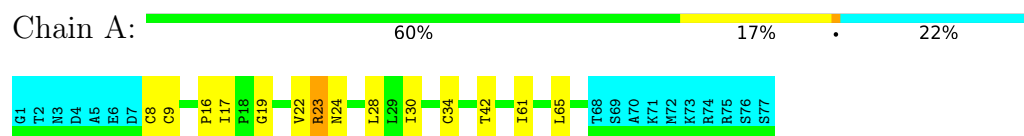
4.2.17 Score per residue for model 17

- Molecule 1: C-C motif chemokine 19



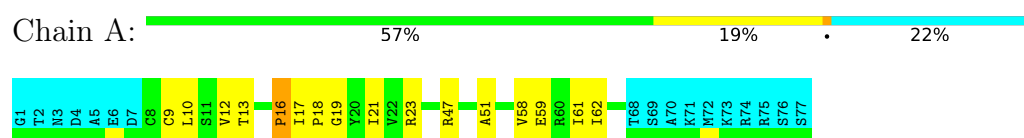
4.2.18 Score per residue for model 18

- Molecule 1: C-C motif chemokine 19



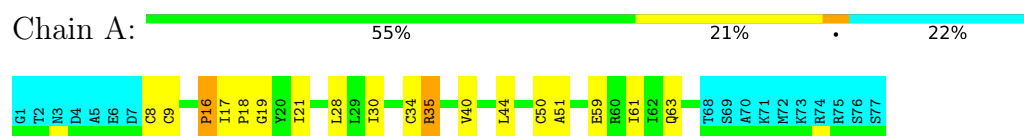
4.2.19 Score per residue for model 19

- Molecule 1: C-C motif chemokine 19



4.2.20 Score per residue for model 20

- Molecule 1: C-C motif chemokine 19



5 Refinement protocol and experimental data overview

The models were refined using the following method: *AUTOMATED METHODS WERE USED FOR BACKBONE CHEMICAL SHIFT ASSIGNMENT AND ITERATIVE NOE REFINEMENT. FINAL STRUCTURES WERE OBTAINED BY MOLECULAR DYNAMICS IN EXPLICIT SOLVENT, AUTOMATED METHODS WERE USED FOR BACKBONE CHEMICAL SHIFT ASSIGNMENT AND ITERATIVE NOE REFINEMENT. FINAL STRUCTURES WERE OBTAINED BY MOLECULAR DYNAMICS IN EXPLICIT SOLVENT.*

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

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

Software name	Classification	Version
Xplor-NIH	refinement	2.9.3
CYANA	refinement	3.1

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	901
Number of shifts mapped to atoms	901
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	83%

6 Model quality i

6.1 Standard geometry i

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

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts i

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	488	509	508	6±2
All	All	9760	10180	10160	128

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:40:VAL:HG12	1:A:50:CYS:SG	0.67	2.30	17	7
1:A:30:ILE:HA	1:A:34:CYS:O	0.61	1.96	20	4
1:A:9:CYS:SG	1:A:40:VAL:HG12	0.58	2.37	20	4
1:A:23:ARG:HD2	1:A:24:ASN:ND2	0.56	2.15	7	1
1:A:58:VAL:O	1:A:62:ILE:HG12	0.55	2.02	13	2
1:A:9:CYS:SG	1:A:38:ALA:HB3	0.54	2.43	12	1
1:A:31:LYS:HE3	1:A:31:LYS:C	0.54	2.23	16	1
1:A:9:CYS:SG	1:A:10:LEU:N	0.50	2.84	2	3
1:A:35:ARG:O	1:A:35:ARG:HD3	0.50	2.07	17	1
1:A:25:PHE:CZ	1:A:62:ILE:HG23	0.49	2.42	16	1
1:A:18:PRO:O	1:A:21:ILE:HG12	0.49	2.07	19	12
1:A:28:LEU:HD21	1:A:34:CYS:SG	0.48	2.48	7	1
1:A:8:CYS:HA	1:A:28:LEU:HG	0.48	1.85	18	2
1:A:23:ARG:HD2	1:A:24:ASN:OD1	0.48	2.08	4	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:59:GLU:O	1:A:63:GLN:HG2	0.48	2.09	1	7
1:A:26:HIS:HB2	1:A:40:VAL:HG22	0.47	1.84	1	1
1:A:22:VAL:HG21	1:A:41:PHE:CD1	0.47	2.45	8	3
1:A:23:ARG:HA	1:A:23:ARG:NE	0.47	2.25	1	1
1:A:35:ARG:HD2	1:A:35:ARG:H	0.46	1.70	20	1
1:A:16:PRO:HG3	1:A:51:ALA:HA	0.46	1.87	8	15
1:A:64:ARG:O	1:A:67:ARG:HG2	0.45	2.11	8	2
1:A:62:ILE:O	1:A:66:GLN:HG3	0.45	2.11	7	2
1:A:58:VAL:O	1:A:62:ILE:HG13	0.45	2.12	3	5
1:A:49:LEU:HD13	1:A:50:CYS:N	0.45	2.27	5	7
1:A:21:ILE:O	1:A:44:LEU:HG	0.45	2.11	9	5
1:A:41:PHE:O	1:A:48:GLN:HA	0.44	2.13	2	3
1:A:67:ARG:HA	1:A:67:ARG:HE	0.44	1.72	15	2
1:A:22:VAL:HG13	1:A:65:LEU:HD11	0.44	1.90	18	7
1:A:8:CYS:SG	1:A:28:LEU:HD12	0.44	2.52	18	2
1:A:19:GLY:HA3	1:A:61:ILE:CD1	0.43	2.43	20	6
1:A:60:ARG:O	1:A:64:ARG:HG3	0.43	2.13	12	1
1:A:10:LEU:HD12	1:A:10:LEU:H	0.43	1.73	19	1
1:A:38:ALA:HA	1:A:53:PRO:HD3	0.43	1.90	14	1
1:A:47:ARG:HE	1:A:47:ARG:HA	0.43	1.73	16	1
1:A:9:CYS:HB2	1:A:40:VAL:HG12	0.42	1.91	16	1
1:A:8:CYS:SG	1:A:9:CYS:N	0.42	2.91	17	2
1:A:31:LYS:HE3	1:A:32:ASP:N	0.42	2.29	16	1
1:A:56:PRO:HA	1:A:59:GLU:OE1	0.42	2.15	5	1
1:A:63:GLN:O	1:A:67:ARG:HG2	0.41	2.16	6	1
1:A:25:PHE:CE1	1:A:65:LEU:HB3	0.41	2.50	11	1
1:A:42:THR:HA	1:A:47:ARG:O	0.41	2.16	16	1
1:A:24:ASN:O	1:A:42:THR:HB	0.40	2.16	18	1
1:A:49:LEU:HD23	1:A:50:CYS:N	0.40	2.32	10	1
1:A:12:VAL:HG12	1:A:50:CYS:CB	0.40	2.47	6	1
1:A:12:VAL:HG23	1:A:13:THR:O	0.40	2.17	19	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	60/77 (78%)	56±1 (93±2%)	3±1 (5±2%)	1±1 (2±1%)	11	52
All	All	1200/1540 (78%)	1120 (93%)	56 (5%)	24 (2%)	11	52

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)
1	A	16	PRO	19
1	A	33	GLY	5

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	55/69 (80%)	52±1 (94±3%)	4±1 (6±3%)	21	70
All	All	1100/1380 (80%)	1030 (94%)	70 (6%)	21	70

All 15 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	17	ILE	20
1	A	23	ARG	15
1	A	59	GLU	8
1	A	9	CYS	8
1	A	12	VAL	4
1	A	31	LYS	2
1	A	67	ARG	2
1	A	47	ARG	2
1	A	49	LEU	2
1	A	35	ARG	2
1	A	48	GLN	1
1	A	63	GLN	1
1	A	16	PRO	1
1	A	28	LEU	1
1	A	64	ARG	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 83% for the well-defined parts and 82% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping (i)

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

Total number of shifts	901
Number of shifts mapped to atoms	901
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

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
$^{13}\text{C}_\alpha$	77	-0.40 ± 0.26	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	73	-0.11 ± 0.13	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	69	0.42 ± 0.57	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 83%, i.e. 739 atoms were assigned a chemical shift out of a possible 889. 0 out of 12 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	229/291 (79%)	116/117 (99%)	60/120 (50%)	53/54 (98%)
Sidechain	466/540 (86%)	317/352 (90%)	143/159 (90%)	6/29 (21%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	44/58 (76%)	26/28 (93%)	17/27 (63%)	1/3 (33%)
Overall	739/889 (83%)	459/497 (92%)	220/306 (72%)	60/86 (70%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 82%, i.e. 901 atoms were assigned a chemical shift out of a possible 1098. 0 out of 12 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	296/377 (79%)	150/152 (99%)	77/154 (50%)	69/71 (97%)
Sidechain	561/663 (85%)	381/429 (89%)	173/196 (88%)	7/38 (18%)
Aromatic	44/58 (76%)	26/28 (93%)	17/27 (63%)	1/3 (33%)
Overall	901/1098 (82%)	557/609 (91%)	267/377 (71%)	77/112 (69%)

7.1.4 Statistically unusual chemical shifts [i](#)

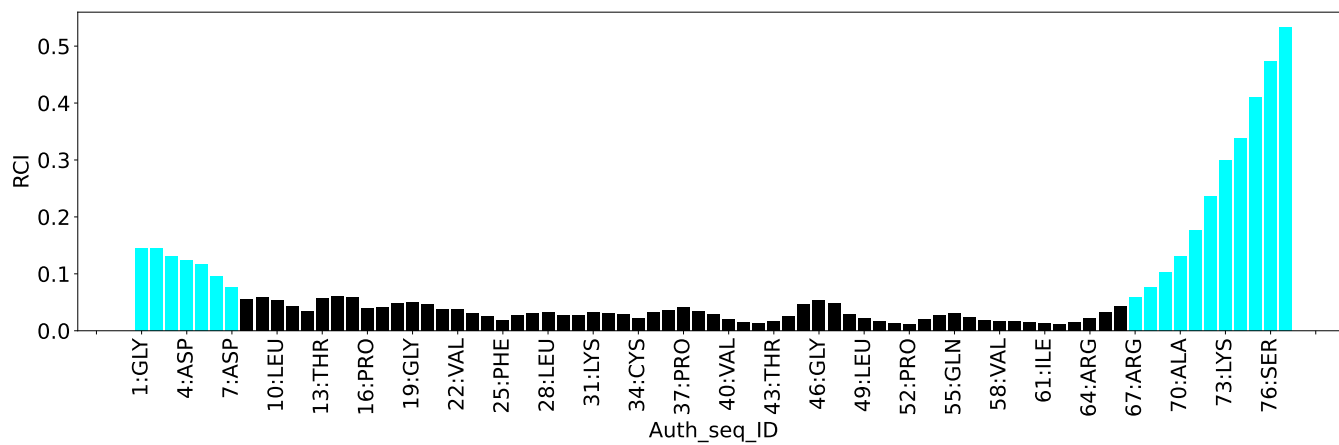
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	52	PRO	HA	1.50	2.78 – 6.00	-9.0

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:



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	1387
Intra-residue ($ i-j =0$)	782
Sequential ($ i-j =1$)	246
Medium range ($ i-j >1$ and $ i-j <5$)	109
Long range ($ i-j \geq 5$)	244
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	6
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	18.0
Number of long range restraints per residue ¹	3.2

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)	5.1	0.2
0.2-0.5 (Medium)	0.3	0.26
>0.5 (Large)	None	None

8.2.2 Average number of dihedral-angle violations per model

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

9 Distance violation analysis [i](#)

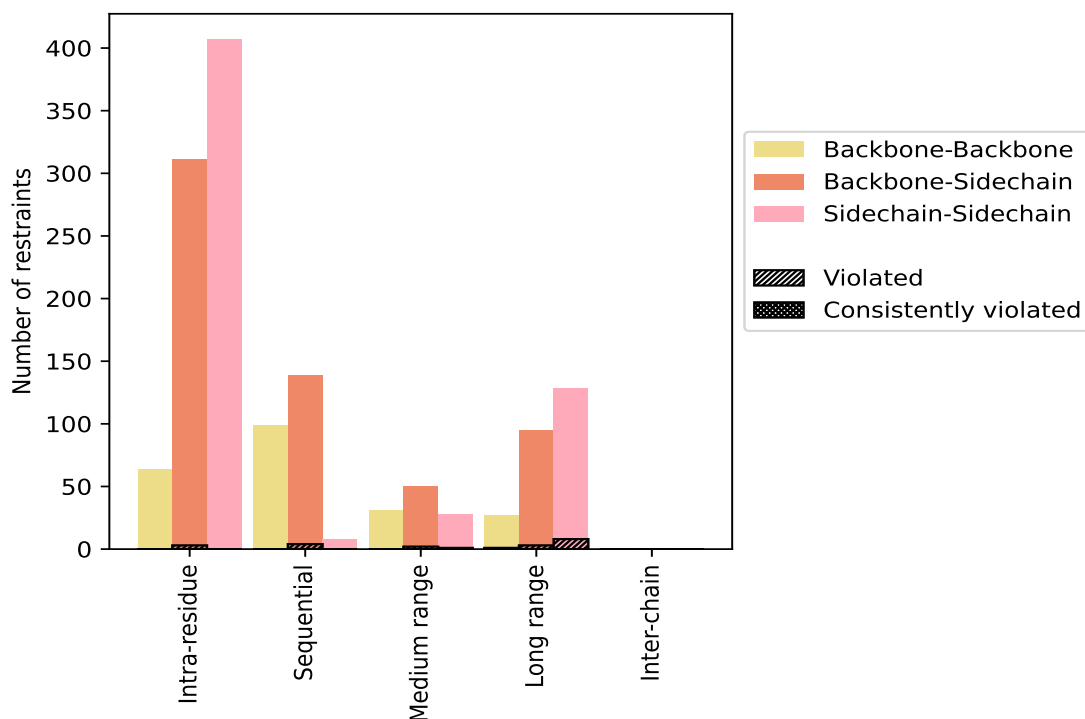
9.1 Summary of distance violations [i](#)

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

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	782	56.4	3	0.4	0.2	0	0.0	0.0
Backbone-Backbone	64	4.6	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	311	22.4	3	1.0	0.2	0	0.0	0.0
Sidechain-Sidechain	407	29.3	0	0.0	0.0	0	0.0	0.0
Sequential ($i-j =1$)	246	17.7	4	1.6	0.3	0	0.0	0.0
Backbone-Backbone	99	7.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	139	10.0	4	2.9	0.3	0	0.0	0.0
Sidechain-Sidechain	8	0.6	0	0.0	0.0	0	0.0	0.0
Medium range ($i-j >1$ & $i-j <5$)	109	7.9	3	2.8	0.2	0	0.0	0.0
Backbone-Backbone	31	2.2	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	50	3.6	2	4.0	0.1	0	0.0	0.0
Sidechain-Sidechain	28	2.0	1	3.6	0.1	0	0.0	0.0
Long range ($i-j \geq 5$)	244	17.6	12	4.9	0.9	0	0.0	0.0
Backbone-Backbone	27	1.9	1	3.7	0.1	0	0.0	0.0
Backbone-Sidechain	95	6.8	3	3.2	0.2	0	0.0	0.0
Sidechain-Sidechain	122	8.8	8	6.6	0.6	0	0.0	0.0
Inter-chain	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Hydrogen bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	6	0.4	0	0.0	0.0	0	0.0	0.0
Total	1387	100.0	22	1.6	1.6	0	0.0	0.0
Backbone-Backbone	221	15.9	1	0.5	0.1	0	0.0	0.0
Backbone-Sidechain	595	42.9	12	2.0	0.9	0	0.0	0.0
Sidechain-Sidechain	571	41.2	9	1.6	0.6	0	0.0	0.0

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

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



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and 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	0	0	0	3	0	3	0.13	0.18	0.03	0.11
2	1	0	2	2	0	5	0.15	0.26	0.05	0.13
3	2	1	0	3	0	6	0.15	0.2	0.04	0.15
4	0	1	0	5	0	6	0.14	0.18	0.03	0.14
5	1	0	1	3	0	5	0.15	0.24	0.05	0.13
6	1	1	0	3	0	5	0.14	0.2	0.03	0.13
7	1	1	0	6	0	8	0.15	0.23	0.04	0.14
8	1	0	1	5	0	7	0.14	0.21	0.04	0.12
9	1	0	1	2	0	4	0.14	0.15	0.01	0.15
10	1	1	0	4	0	6	0.16	0.25	0.05	0.14
11	0	0	0	3	0	3	0.15	0.17	0.02	0.16

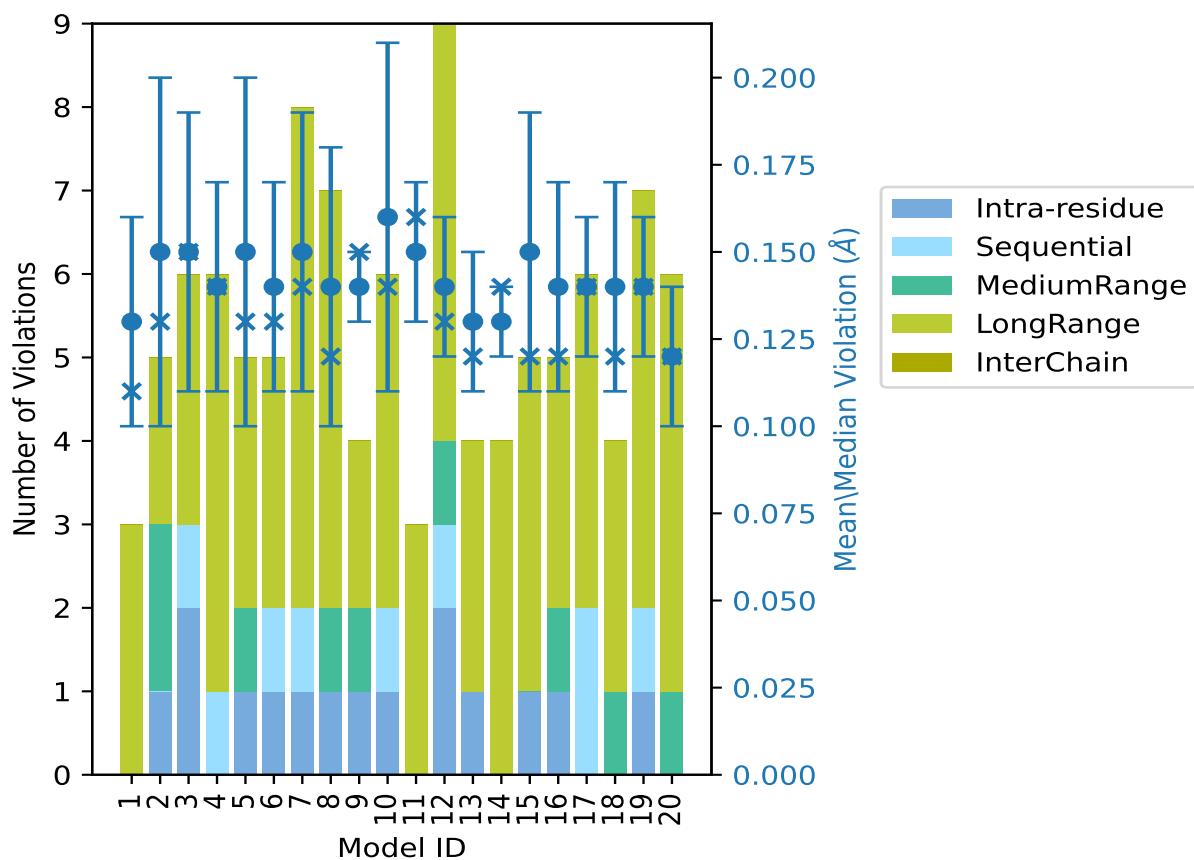
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
12	2	1	1	5	0	9	0.14	0.17	0.02	0.13
13	1	0	0	3	0	4	0.13	0.16	0.02	0.12
14	0	0	0	4	0	4	0.13	0.14	0.01	0.14
15	1	0	0	4	0	5	0.15	0.22	0.04	0.12
16	1	0	1	3	0	5	0.14	0.18	0.03	0.12
17	0	2	0	4	0	6	0.14	0.17	0.02	0.14
18	0	0	1	3	0	4	0.14	0.19	0.03	0.12
19	1	1	0	5	0	7	0.14	0.16	0.02	0.14
20	0	0	1	5	0	6	0.12	0.18	0.02	0.12

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

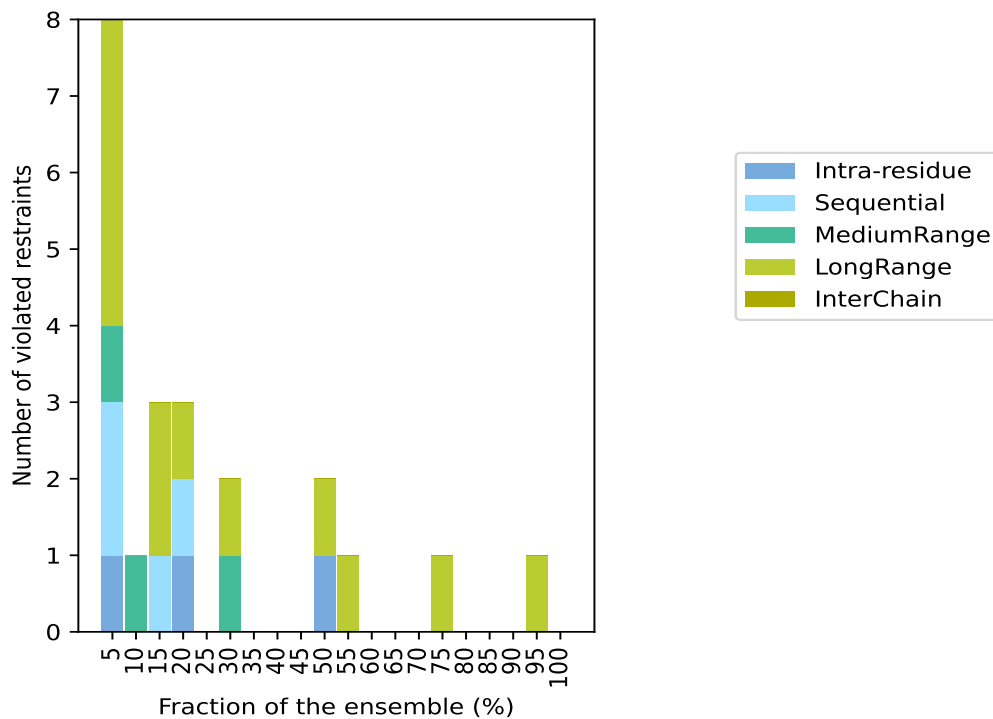
9.3 Distance violation statistics for the ensemble

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, 1359(IR:779, SQ:242, MR:106, LR:232, 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	%
1	2	1	4	0	8	1	5.0
0	0	1	0	0	1	2	10.0
0	1	0	2	0	3	3	15.0
1	1	0	1	0	3	4	20.0
0	0	0	0	0	0	5	25.0
0	0	1	1	0	2	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	1	0	2	10	50.0
0	0	0	1	0	1	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	1	0	1	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	1	0	1	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

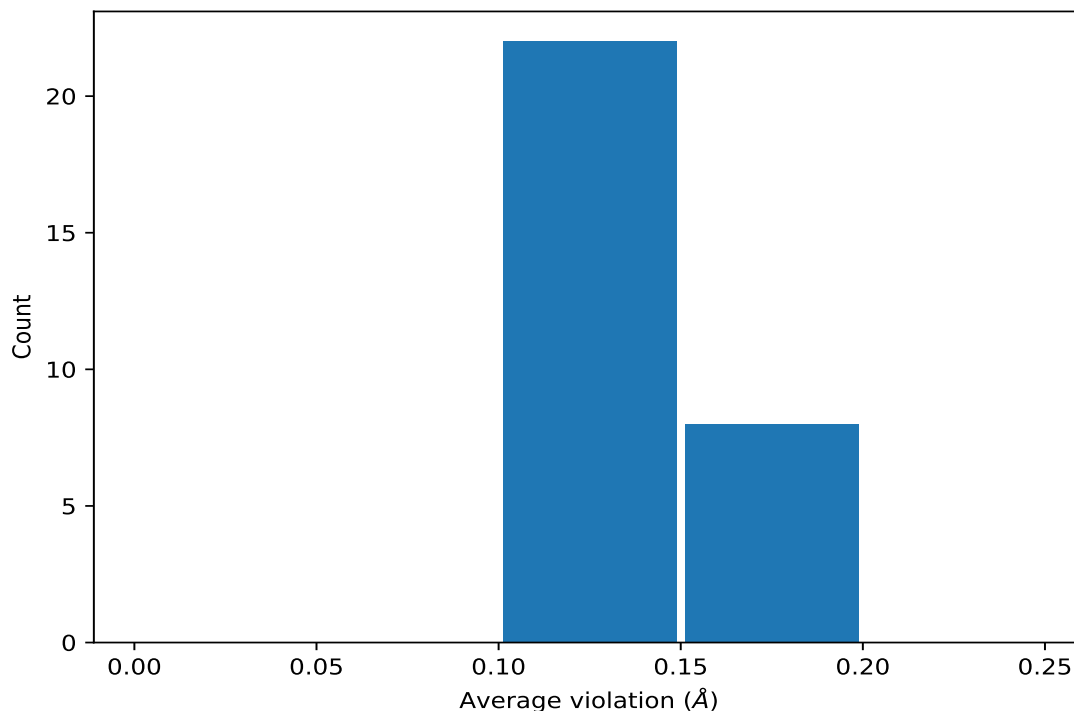
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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)
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	19	0.17	0.03	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	19	0.17	0.03	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	19	0.17	0.03	0.17
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	15	0.12	0.01	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	15	0.12	0.01	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	15	0.12	0.01	0.11
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	11	0.18	0.05	0.18
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	10	0.13	0.02	0.12
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	10	0.13	0.02	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	10	0.13	0.01	0.13
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	10	0.13	0.01	0.13
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	10	0.13	0.01	0.13
(1,607)	1:A:30:ILE:HG21	1:A:33:GLY:HA3	6	0.14	0.02	0.14
(1,607)	1:A:30:ILE:HG22	1:A:33:GLY:HA3	6	0.14	0.02	0.14
(1,607)	1:A:30:ILE:HG23	1:A:33:GLY:HA3	6	0.14	0.02	0.14
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG21	6	0.13	0.02	0.12

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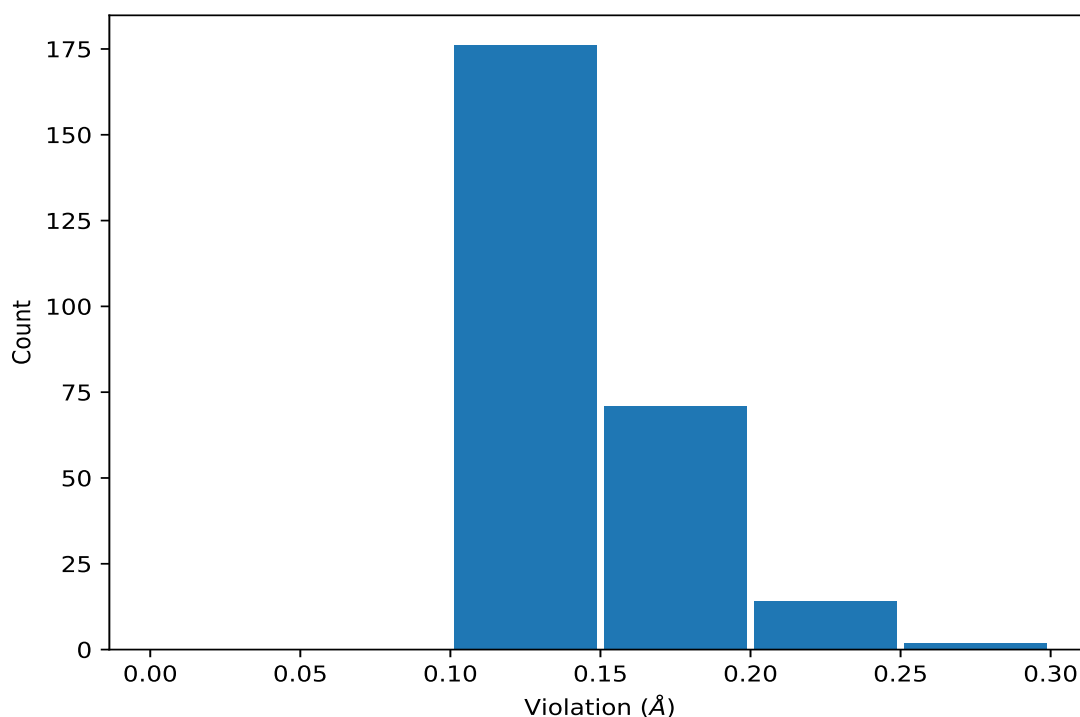
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG22	6	0.13	0.02	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG23	6	0.13	0.02	0.12
(1,940)	1:A:49:LEU:HD11	1:A:50:CYS:H	4	0.16	0.01	0.16
(1,940)	1:A:49:LEU:HD12	1:A:50:CYS:H	4	0.16	0.01	0.16
(1,940)	1:A:49:LEU:HD13	1:A:50:CYS:H	4	0.16	0.01	0.16
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG21	4	0.12	0.01	0.12
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG22	4	0.12	0.01	0.12
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG23	4	0.12	0.01	0.12
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD1	4	0.12	0.01	0.11
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD2	4	0.12	0.01	0.11
(1,935)	1:A:49:LEU:HG	1:A:50:CYS:H	3	0.16	0.01	0.16
(1,295)	1:A:20:TYR:H	1:A:61:ILE:HB	3	0.12	0.01	0.13
(1,789)	1:A:41:PHE:HZ	1:A:61:ILE:HB	3	0.11	0.0	0.11
(1,566)	1:A:29:LEU:HG	1:A:32:ASP:H	2	0.12	0.0	0.12

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)
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	2	0.26
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	10	0.25
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	5	0.24
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	5	0.24
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	5	0.24
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	7	0.23
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	15	0.22
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	15	0.22
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	15	0.22
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	8	0.21
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	8	0.21
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	8	0.21
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	6	0.2
(1,1257)	1:A:65:LEU:H	1:A:65:LEU:HD11	3	0.2
(1,1257)	1:A:65:LEU:H	1:A:65:LEU:HD12	3	0.2
(1,1257)	1:A:65:LEU:H	1:A:65:LEU:HD13	3	0.2
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	3	0.19
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	3	0.19
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	3	0.19
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	18	0.19
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	18	0.19
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	18	0.19
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	4	0.18
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	20	0.18
(1,607)	1:A:30:ILE:HG21	1:A:33:GLY:HA3	8	0.18
(1,607)	1:A:30:ILE:HG22	1:A:33:GLY:HA3	8	0.18
(1,607)	1:A:30:ILE:HG23	1:A:33:GLY:HA3	8	0.18
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	16	0.18
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	16	0.18
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	1	0.18
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	1	0.18
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	1	0.18
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	4	0.18
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	4	0.18
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	4	0.18
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	16	0.18
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	16	0.18

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	16	0.18
(1,940)	1:A:49:LEU:HD11	1:A:50:CYS:H	3	0.17
(1,940)	1:A:49:LEU:HD12	1:A:50:CYS:H	3	0.17
(1,940)	1:A:49:LEU:HD13	1:A:50:CYS:H	3	0.17
(1,940)	1:A:49:LEU:HD11	1:A:50:CYS:H	10	0.17
(1,940)	1:A:49:LEU:HD12	1:A:50:CYS:H	10	0.17
(1,940)	1:A:49:LEU:HD13	1:A:50:CYS:H	10	0.17
(1,935)	1:A:49:LEU:HG	1:A:50:CYS:H	17	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	11	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	11	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	11	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	12	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	12	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	12	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	17	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	17	0.17
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	17	0.17
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	19	0.16
(1,935)	1:A:49:LEU:HG	1:A:50:CYS:H	12	0.16
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG21	11	0.16
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG22	11	0.16
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG23	11	0.16
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	15	0.16
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	15	0.16
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	15	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	7	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	7	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	7	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	13	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	13	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	13	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	19	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	19	0.16
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	19	0.16
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	17	0.15
(1,940)	1:A:49:LEU:HD11	1:A:50:CYS:H	4	0.15
(1,940)	1:A:49:LEU:HD12	1:A:50:CYS:H	4	0.15
(1,940)	1:A:49:LEU:HD13	1:A:50:CYS:H	4	0.15
(1,935)	1:A:49:LEU:HG	1:A:50:CYS:H	7	0.15
(1,607)	1:A:30:ILE:HG21	1:A:33:GLY:HA3	2	0.15
(1,607)	1:A:30:ILE:HG22	1:A:33:GLY:HA3	2	0.15
(1,607)	1:A:30:ILE:HG23	1:A:33:GLY:HA3	2	0.15

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,607)	1:A:30:ILE:HG21	1:A:33:GLY:HA3	9	0.15
(1,607)	1:A:30:ILE:HG22	1:A:33:GLY:HA3	9	0.15
(1,607)	1:A:30:ILE:HG23	1:A:33:GLY:HA3	9	0.15
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	19	0.15
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	19	0.15
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	9	0.15
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	9	0.15
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	9	0.15
(1,940)	1:A:49:LEU:HD11	1:A:50:CYS:H	19	0.14
(1,940)	1:A:49:LEU:HD12	1:A:50:CYS:H	19	0.14
(1,940)	1:A:49:LEU:HD13	1:A:50:CYS:H	19	0.14
(1,548)	1:A:28:LEU:HB3	1:A:40:VAL:HG11	14	0.14
(1,548)	1:A:28:LEU:HB3	1:A:40:VAL:HG12	14	0.14
(1,548)	1:A:28:LEU:HB3	1:A:40:VAL:HG13	14	0.14
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	8	0.14
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	8	0.14
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	9	0.14
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	9	0.14
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	7	0.14
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	7	0.14
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	7	0.14
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	12	0.14
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	12	0.14
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	12	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	10	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	10	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	10	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	12	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	12	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	12	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	14	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	14	0.14
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	14	0.14
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	10	0.14
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	10	0.14
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	10	0.14
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	12	0.13
(1,607)	1:A:30:ILE:HG21	1:A:33:GLY:HA3	5	0.13
(1,607)	1:A:30:ILE:HG22	1:A:33:GLY:HA3	5	0.13
(1,607)	1:A:30:ILE:HG23	1:A:33:GLY:HA3	5	0.13
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	5	0.13
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	5	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG21	7	0.13
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG22	7	0.13
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG23	7	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	5	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	5	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	5	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	6	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	6	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	6	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	10	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	10	0.13
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	10	0.13
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG21	17	0.13
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG22	17	0.13
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG23	17	0.13
(1,295)	1:A:20:TYR:H	1:A:61:ILE:HB	4	0.13
(1,295)	1:A:20:TYR:H	1:A:61:ILE:HB	13	0.13
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD1	12	0.13
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD2	12	0.13
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	6	0.13
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	6	0.13
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	6	0.13
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	17	0.13
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	17	0.13
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	17	0.13
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	2	0.13
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	2	0.13
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	2	0.13
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	14	0.13
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	14	0.13
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	14	0.13
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	15	0.12
(1,607)	1:A:30:ILE:HG21	1:A:33:GLY:HA3	20	0.12
(1,607)	1:A:30:ILE:HG22	1:A:33:GLY:HA3	20	0.12
(1,607)	1:A:30:ILE:HG23	1:A:33:GLY:HA3	20	0.12
(1,566)	1:A:29:LEU:HG	1:A:32:ASP:H	16	0.12
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	2	0.12
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	2	0.12
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	12	0.12
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	12	0.12
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	13	0.12
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	13	0.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	15	0.12
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	15	0.12
(1,431)	1:A:24:ASN:HB2	1:A:42:THR:HG21	12	0.12
(1,431)	1:A:24:ASN:HB2	1:A:42:THR:HG22	12	0.12
(1,431)	1:A:24:ASN:HB2	1:A:42:THR:HG23	12	0.12
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG21	18	0.12
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG22	18	0.12
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG23	18	0.12
(1,40)	1:A:7:ASP:HA	1:A:28:LEU:HD11	11	0.12
(1,40)	1:A:7:ASP:HA	1:A:28:LEU:HD12	11	0.12
(1,40)	1:A:7:ASP:HA	1:A:28:LEU:HD13	11	0.12
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	9	0.12
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	9	0.12
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	9	0.12
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	16	0.12
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	16	0.12
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	16	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG21	7	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG22	7	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG23	7	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG21	8	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG22	8	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG23	8	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG21	19	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG22	19	0.12
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG23	19	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	3	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	3	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	3	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	5	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	5	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	5	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	7	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	7	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	7	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG21	20	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG22	20	0.12
(1,208)	1:A:16:PRO:HG3	1:A:58:VAL:HG23	20	0.12
(1,114)	1:A:12:VAL:HB	1:A:16:PRO:HB2	18	0.12
(1,99)	1:A:11:SER:H	1:A:50:CYS:H	14	0.11
(1,789)	1:A:41:PHE:HZ	1:A:61:ILE:HB	4	0.11
(1,789)	1:A:41:PHE:HZ	1:A:61:ILE:HB	8	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,789)	1:A:41:PHE:HZ	1:A:61:ILE:HB	16	0.11
(1,69)	1:A:9:CYS:HB2	1:A:28:LEU:HD11	20	0.11
(1,69)	1:A:9:CYS:HB2	1:A:28:LEU:HD12	20	0.11
(1,69)	1:A:9:CYS:HB2	1:A:28:LEU:HD13	20	0.11
(1,69)	1:A:9:CYS:HB3	1:A:28:LEU:HD11	20	0.11
(1,69)	1:A:9:CYS:HB3	1:A:28:LEU:HD12	20	0.11
(1,69)	1:A:9:CYS:HB3	1:A:28:LEU:HD13	20	0.11
(1,607)	1:A:30:ILE:HG21	1:A:33:GLY:HA3	12	0.11
(1,607)	1:A:30:ILE:HG22	1:A:33:GLY:HA3	12	0.11
(1,607)	1:A:30:ILE:HG23	1:A:33:GLY:HA3	12	0.11
(1,566)	1:A:29:LEU:HG	1:A:32:ASP:H	2	0.11
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE1	6	0.11
(1,502)	1:A:27:TYR:HA	1:A:27:TYR:HE2	6	0.11
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG21	8	0.11
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG22	8	0.11
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG23	8	0.11
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG21	19	0.11
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG22	19	0.11
(1,401)	1:A:23:ARG:HB2	1:A:42:THR:HG23	19	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	1	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	1	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	1	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	3	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	3	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	3	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	4	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	4	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	4	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	13	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	13	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	13	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	15	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	15	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	15	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	18	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	18	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	18	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	19	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	19	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	19	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD11	20	0.11
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD12	20	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,355)	1:A:22:VAL:HA	1:A:65:LEU:HD13	20	0.11
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG21	1	0.11
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG22	1	0.11
(1,307)	1:A:20:TYR:HB3	1:A:61:ILE:HG23	1	0.11
(1,295)	1:A:20:TYR:H	1:A:61:ILE:HB	8	0.11
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD1	3	0.11
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD2	3	0.11
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD1	7	0.11
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD2	7	0.11
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD1	10	0.11
(1,293)	1:A:20:TYR:H	1:A:20:TYR:HD2	10	0.11
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB1	20	0.11
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB2	20	0.11
(1,195)	1:A:16:PRO:HD3	1:A:51:ALA:HB3	20	0.11
(1,17)	1:A:4:ASP:HB3	1:A:5:ALA:H	17	0.11
(1,1343)	1:A:71:LYS:HB2	1:A:72:MET:H	6	0.11

10 Dihedral-angle violation analysis

Dihedral angle analysis failed due to data error in the dihedral angle restraints, possibly missing target value