wwPDB NMR Structure Validation Summary Report (i)

Jun 4, 2023 – 10:46 PM EDT

PDB ID	:	2LZU
BMRB ID	:	18778
Title	:	Solution structure of LIMD2
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Deposited on	:	2012-10-10

This is a wwPDB NMR Structure Validation Summary 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 (i) 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 (1)) were used in the production of this report:

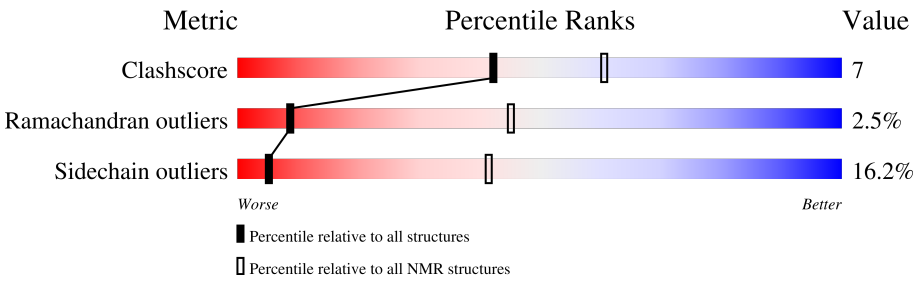
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
BMRB Restraints Analysis	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
SOLUTION NMR

The overall completeness of chemical shifts assignment is 73%.

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	72	56% 24% • 18%

2 Ensemble composition and analysis (i)

This entry contains 10 models. Model 10 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:41-A:99 (59)	0.43	10

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

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

The models can be grouped into 3 clusters. No single-model clusters were found.

Cluster number	Models
1	1, 7, 8, 10
2	2, 5, 9
3	3, 4, 6

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1144 atoms, of which 562 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called LIM domain-containing protein 2.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	72	1142	368	562	102	102	8	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	
			Total	Zn
2	A	2	2	2

4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

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

• Molecule 1: LIM domain-containing protein 2

Chain A: 56% 24% • 18%

R36 A37 Q38 V39 K40 E41 T42 C43 Q47 K48 T49 V50 D59 I62 F63 H64 C67 C70 K71 H72 C73 K76 L77 S78 L79 Y82 Y90 Q96 K100 S101 K102 G103 N104 Y105 D106 E107

4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 10. Colouring as in section 4.1 above.

• Molecule 1: LIM domain-containing protein 2

Chain A: 54% 25% • 18%

R36 A37 Q38 V39 K40 E41 T42 C43 Q47 K48 T49 V50 D59 I62 F63 H64 C67 C70 K71 H72 C73 K76 L77 S78 L79 Y82 Y90 Q96 K100 S101 K102 G103 N104 Y105 D106 E107

5 Refinement protocol and experimental data overview (i)

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

Of the 450 calculated structures, 10 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
ARIA	structure solution	2.3
CNS	refinement	

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	3
Total number of shifts	1396
Number of shifts mapped to atoms	1396
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	73%

6 Model quality (i)

6.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	473	453	451	7±1
All	All	4750	4530	4510	67

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

5 of 15 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:43:CYS:O	1:A:47:GLN:HA	0.67	1.89	2	10
1:A:64:HIS:HB2	1:A:67:CYS:SG	0.66	2.30	1	10
1:A:70:CYS:HB3	1:A:73:CYS:O	0.55	2.02	8	9
1:A:43:CYS:HA	1:A:62:ILE:O	0.55	2.01	5	9
1:A:43:CYS:HB3	1:A:48:LYS:O	0.53	2.03	5	10

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	59/72 (82%)	54±1 (91±2%)	4±1 (7±2%)	2±1 (3±2%)	9 45
All	All	590/720 (82%)	535 (91%)	40 (7%)	15 (3%)	9 45

All 4 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	79	LEU	6
1	A	80	GLY	5
1	A	59	ASP	3
1	A	60	LYS	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	52/63 (83%)	44±2 (84±3%)	8±2 (16±3%)	5 41
All	All	520/630 (83%)	436 (84%)	84 (16%)	5 41

5 of 12 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	42	THR	10
1	A	43	CYS	10
1	A	63	PHE	10
1	A	79	LEU	10
1	A	90	TYR	10

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)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 73% for the well-defined parts and 63% 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	305
Number of shifts mapped to atoms	305
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing (i)

No chemical shift referencing corrections were calculated (not enough data).

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 38%, i.e. 305 atoms were assigned a chemical shift out of a possible 798. 0 out of 8 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	**^{1}H**	**^{13}C**	**^{15}N**
Backbone	55/293 (19%)	55/118 (47%)	0/118 (0%)	0/57 (0%)
Sidechain	205/391 (52%)	205/256 (80%)	0/123 (0%)	0/12 (0%)
Aromatic	45/114 (39%)	45/57 (79%)	0/50 (0%)	0/7 (0%)
Overall	305/798 (38%)	305/431 (71%)	0/291 (0%)	0/76 (0%)

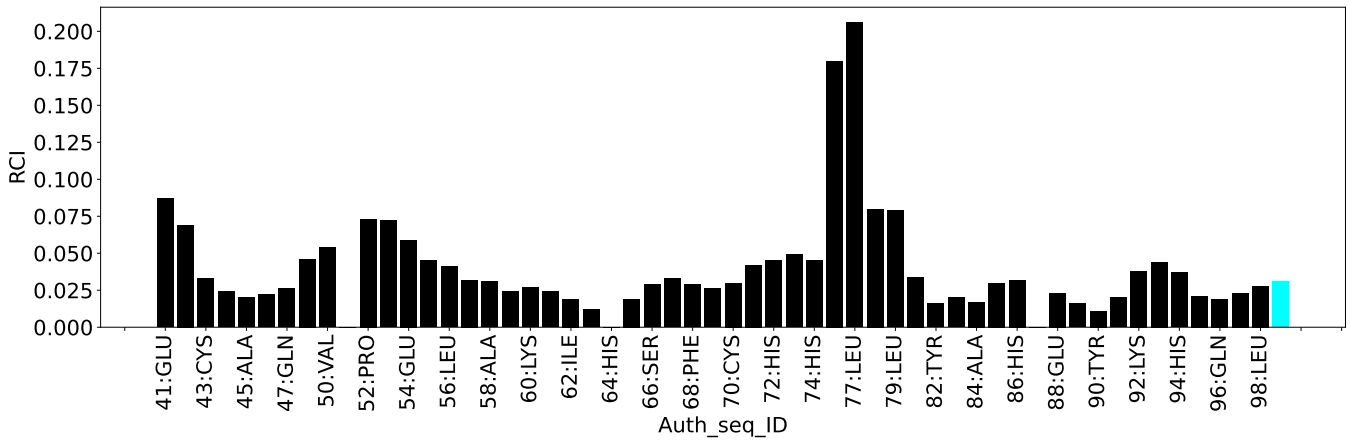
7.1.4 Statistically unusual chemical shifts (i)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots (i)

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

Random coil index (RCI) for chain A:



7.2 Chemical shift list 2

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_2*

7.2.1 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	617
Number of shifts mapped to atoms	617
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	76

7.2.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	51	1.64 ± 0.28	Should be applied
$^{13}C_{\beta}$	47	1.37 ± 0.46	Should be applied
$^{13}C'$	0	—	None (insufficient data)
^{15}N	60	0.16 ± 0.62	None needed (< 0.5 ppm)

7.2.3 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 73%, i.e. 581 atoms were assigned a chemical shift out of a possible 798. 0 out of 8 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	218/293 (74%)	113/118 (96%)	51/118 (43%)	54/57 (95%)
Sidechain	318/391 (81%)	233/256 (91%)	85/123 (69%)	0/12 (0%)
Aromatic	45/114 (39%)	45/57 (79%)	0/50 (0%)	0/7 (0%)
Overall	581/798 (73%)	391/431 (91%)	136/291 (47%)	54/76 (71%)

7.2.4 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
2	A	93	PRO	CG	52.90	21.69 – 32.72	23.3
2	A	49	THR	CG2	46.81	16.06 – 27.03	23.0
2	A	56	LEU	CG	51.61	21.37 – 32.19	22.9
2	A	61	LEU	CG	51.58	21.37 – 32.19	22.9
2	A	52	PRO	CG	52.28	21.69 – 32.72	22.7
2	A	98	LEU	CG	51.31	21.37 – 32.19	22.7
2	A	48	LYS	CG	49.84	19.35 – 30.45	22.5
2	A	85	LEU	CG	50.71	21.37 – 32.19	22.1
2	A	48	LYS	CD	53.09	23.50 – 34.42	22.1
2	A	42	THR	CG2	45.75	16.06 – 27.03	22.1
2	A	77	LEU	CG	49.00	21.37 – 32.19	20.5
2	A	62	ILE	CG2	42.83	10.93 – 24.12	19.2
2	A	50	VAL	CG1	45.31	14.71 – 28.29	17.5
2	A	57	VAL	CG1	45.21	14.71 – 28.29	17.5
2	A	78	SER	CB	38.82	56.28 – 71.32	-16.6
2	A	56	LEU	CD1	50.80	16.71 – 32.55	16.5
2	A	61	LEU	CD1	50.77	16.71 – 32.55	16.5

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List Id	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
2	A	93	PRO	CA	39.06	55.85 – 70.84	-16.2
2	A	49	THR	CB	41.94	61.12 – 78.27	-16.2
2	A	84	ALA	CB	47.32	10.19 – 27.75	16.1
2	A	52	PRO	CA	39.32	55.85 – 70.84	-16.0
2	A	75	THR	CB	42.23	61.12 – 78.27	-16.0
2	A	57	VAL	CG2	45.21	13.71 – 28.88	15.8
2	A	79	LEU	CD1	49.44	16.71 – 32.55	15.7
2	A	77	LEU	CD1	49.40	16.71 – 32.55	15.7
2	A	50	VAL	CG2	44.90	13.71 – 28.88	15.6
2	A	85	LEU	CD2	49.90	15.73 – 32.47	15.4
2	A	98	LEU	CD1	48.88	16.71 – 32.55	15.3
2	A	56	LEU	CD2	49.62	15.73 – 32.47	15.2
2	A	42	THR	CB	43.78	61.12 – 78.27	-15.1
2	A	45	ALA	CB	45.11	10.19 – 27.75	14.9
2	A	62	ILE	CG1	52.76	19.24 – 36.26	14.7
2	A	85	LEU	CD1	47.87	16.71 – 32.55	14.7
2	A	62	ILE	CD1	37.00	5.18 – 21.60	14.4
2	A	79	LEU	CD2	47.81	15.73 – 32.47	14.2
2	A	61	LEU	CD2	47.53	15.73 – 32.47	14.0
2	A	97	GLN	CB	53.75	20.34 – 37.98	13.9
2	A	44	ALA	CB	43.08	10.19 – 27.75	13.7
2	A	98	LEU	CD2	46.84	15.73 – 32.47	13.6
2	A	77	LEU	CD2	46.79	15.73 – 32.47	13.6
2	A	60	LYS	CB	55.94	24.03 – 41.47	13.3
2	A	96	GLN	CB	52.53	20.34 – 37.98	13.2
2	A	66	SER	CA	32.63	48.46 – 68.96	-12.7
2	A	65	ASN	CA	30.60	44.28 – 62.79	-12.4
2	A	97	GLN	CA	30.87	46.17 – 66.97	-12.4
2	A	76	LYS	CA	30.35	46.18 – 67.77	-12.3
2	A	47	GLN	CA	31.16	46.17 – 66.97	-12.2
2	A	96	GLN	CA	31.20	46.17 – 66.97	-12.2
2	A	92	LYS	CB	53.91	24.03 – 41.47	12.1
2	A	60	LYS	CA	31.16	46.18 – 67.77	-12.0
2	A	71	LYS	CA	31.16	46.18 – 67.77	-12.0
2	A	47	GLN	CB	50.14	20.34 – 37.98	11.9
2	A	53	MET	CA	30.35	45.26 – 67.07	-11.8
2	A	55	ARG	CA	30.35	45.44 – 68.13	-11.7
2	A	42	THR	CA	32.75	49.41 – 75.05	-11.5
2	A	94	HIS	CB	53.88	19.76 – 40.75	11.2
2	A	62	ILE	CA	31.97	48.30 – 75.08	-11.1
2	A	72	HIS	CA	31.41	45.04 – 67.94	-10.9
2	A	53	MET	CB	55.94	22.22 – 43.61	10.8

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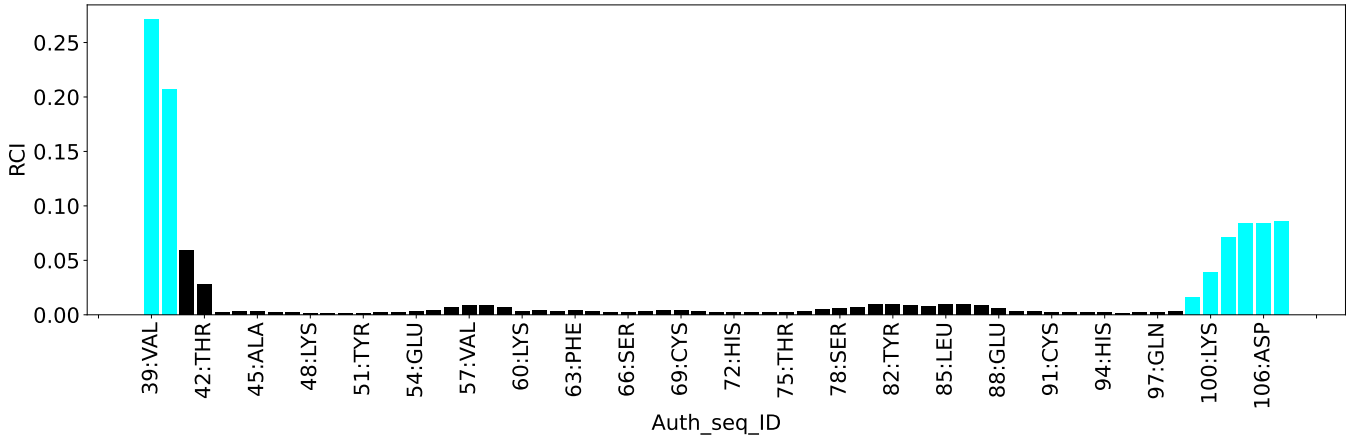
List Id	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
2	A	99	PHE	CA	30.75	45.38 – 70.89	-10.7
2	A	86	HIS	CB	52.74	19.76 – 40.75	10.7
2	A	89	PHE	CA	30.90	45.38 – 70.89	-10.7
2	A	50	VAL	CA	33.13	48.38 – 76.73	-10.4
2	A	74	HIS	CB	51.88	19.76 – 40.75	10.3
2	A	75	THR	CA	35.89	49.41 – 75.05	-10.3
2	A	92	LYS	CA	34.89	46.18 – 67.77	-10.2
2	A	94	HIS	CA	33.30	45.04 – 67.94	-10.1
2	A	49	THR	CA	36.94	49.41 – 75.05	-9.9
2	A	95	PHE	CA	34.12	45.38 – 70.89	-9.4
2	A	43	CYS	CA	30.35	40.80 – 75.33	-8.0
2	A	91	CYS	CA	30.35	40.80 – 75.33	-8.0
2	A	73	CYS	CA	30.75	40.80 – 75.33	-7.9
2	A	70	CYS	CA	31.16	40.80 – 75.33	-7.8
2	A	46	CYS	CA	31.38	40.80 – 75.33	-7.7
2	A	67	CYS	CA	33.44	40.80 – 75.33	-7.1
2	A	87	GLY	N	128.59	91.59 – 127.52	5.3

7.2.5 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:



7.3 Chemical shift list 3

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_3*

7.3.1 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	474
Number of shifts mapped to atoms	474
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

7.3.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	0	—	None (insufficient data)
$^{13}C_{\beta}$	0	—	None (insufficient data)
$^{13}C'$	0	—	None (insufficient data)
^{15}N	60	0.08 ± 1.67	None needed (< 0.5 ppm)

7.3.3 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 55%, i.e. 438 atoms were assigned a chemical shift out of a possible 798. 0 out of 8 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	166/293 (57%)	112/118 (95%)	0/118 (0%)	54/57 (95%)
Sidechain	227/391 (58%)	227/256 (89%)	0/123 (0%)	0/12 (0%)
Aromatic	45/114 (39%)	45/57 (79%)	0/50 (0%)	0/7 (0%)
Overall	438/798 (55%)	384/431 (89%)	0/291 (0%)	54/76 (71%)

7.3.4 Statistically unusual chemical shifts (i)

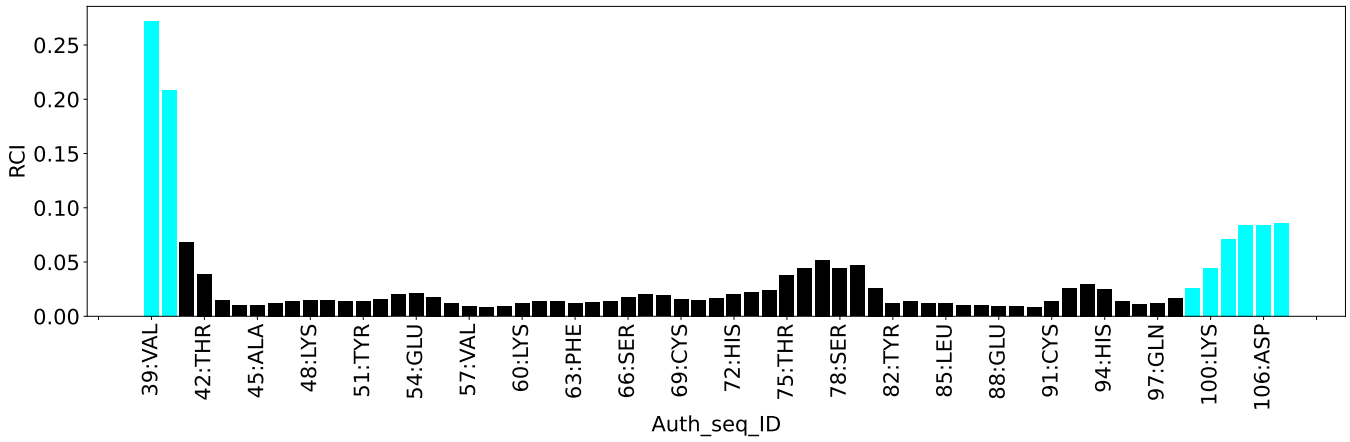
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
3	A	87	GLY	N	128.59	91.59 – 127.52	5.3

7.3.5 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 (i)

8.1 Conformationally restricting restraints (i)

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	1179
Intra-residue (|i-j|=0)	549
Sequential (|i-j|=1)	211
Medium range (|i-j|>1 and |i-j|<5)	124
Long range (|i-j|≥5)	271
Inter-chain	0
Hydrogen bond restraints	24
Disulfide bond restraints	0
Total dihedral-angle restraints	126
Number of unmapped restraints	0
Number of restraints per residue	18.1
Number of long range restraints per residue[1]	3.9

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

8.2 Residual restraint violations (i)

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 (i)

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)	17.0	0.2
0.2-0.5 (Medium)	1.0	0.25
>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)	25.3	4.9
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

9 Distance violation analysis (i)

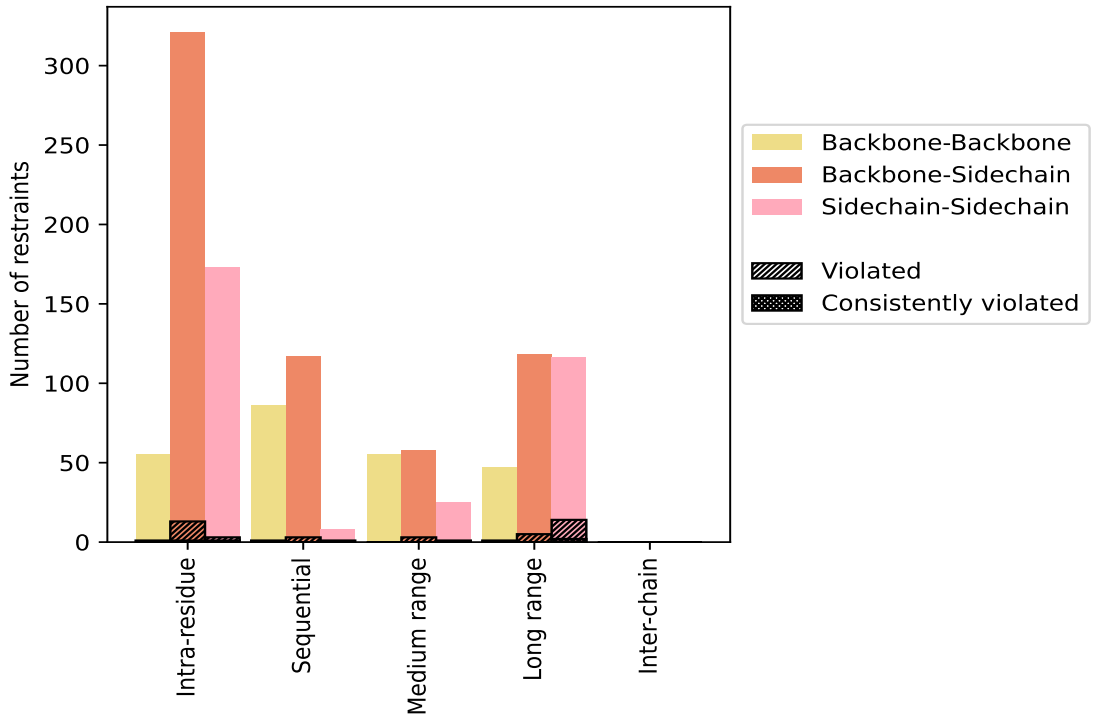
9.1 Summary of distance violations (i)

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.

Restraints type	Count	%[1]	Violated[3]			Consistently Violated[4]		
			Count	%[2]	%[1]	Count	%[2]	%[1]
Intra-residue (|i-j|=0)	549	46.6	17	3.1	1.4	2	0.4	0.2
Backbone-Backbone	55	4.7	1	1.8	0.1	0	0.0	0.0
Backbone-Sidechain	321	27.2	13	4.0	1.1	1	0.3	0.1
Sidechain-Sidechain	173	14.7	3	1.7	0.3	1	0.6	0.1
Sequential (|i-j|=1)	211	17.9	5	2.4	0.4	0	0.0	0.0
Backbone-Backbone	86	7.3	1	1.2	0.1	0	0.0	0.0
Backbone-Sidechain	117	9.9	3	2.6	0.3	0	0.0	0.0
Sidechain-Sidechain	8	0.7	1	12.5	0.1	0	0.0	0.0
Medium range (|i-j|>1 & |i-j|<5)	124	10.5	4	3.2	0.3	0	0.0	0.0
Backbone-Backbone	41	3.5	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	58	4.9	3	5.2	0.3	0	0.0	0.0
Sidechain-Sidechain	25	2.1	1	4.0	0.1	0	0.0	0.0
Long range (|i-j|≥5)	271	23.0	20	7.4	1.7	2	0.7	0.2
Backbone-Backbone	37	3.1	1	2.7	0.1	0	0.0	0.0
Backbone-Sidechain	118	10.0	5	4.2	0.4	0	0.0	0.0
Sidechain-Sidechain	116	9.8	14	12.1	1.2	2	1.7	0.2
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	24	2.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	1179	100.0	46	3.9	3.9	4	0.3	0.3
Backbone-Backbone	243	20.6	3	1.2	0.3	0	0.0	0.0
Backbone-Sidechain	614	52.1	24	3.9	2.0	1	0.2	0.1
Sidechain-Sidechain	322	27.3	19	5.9	1.6	3	0.9	0.3

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

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



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

9.2 Distance violation statistics for each model (i)

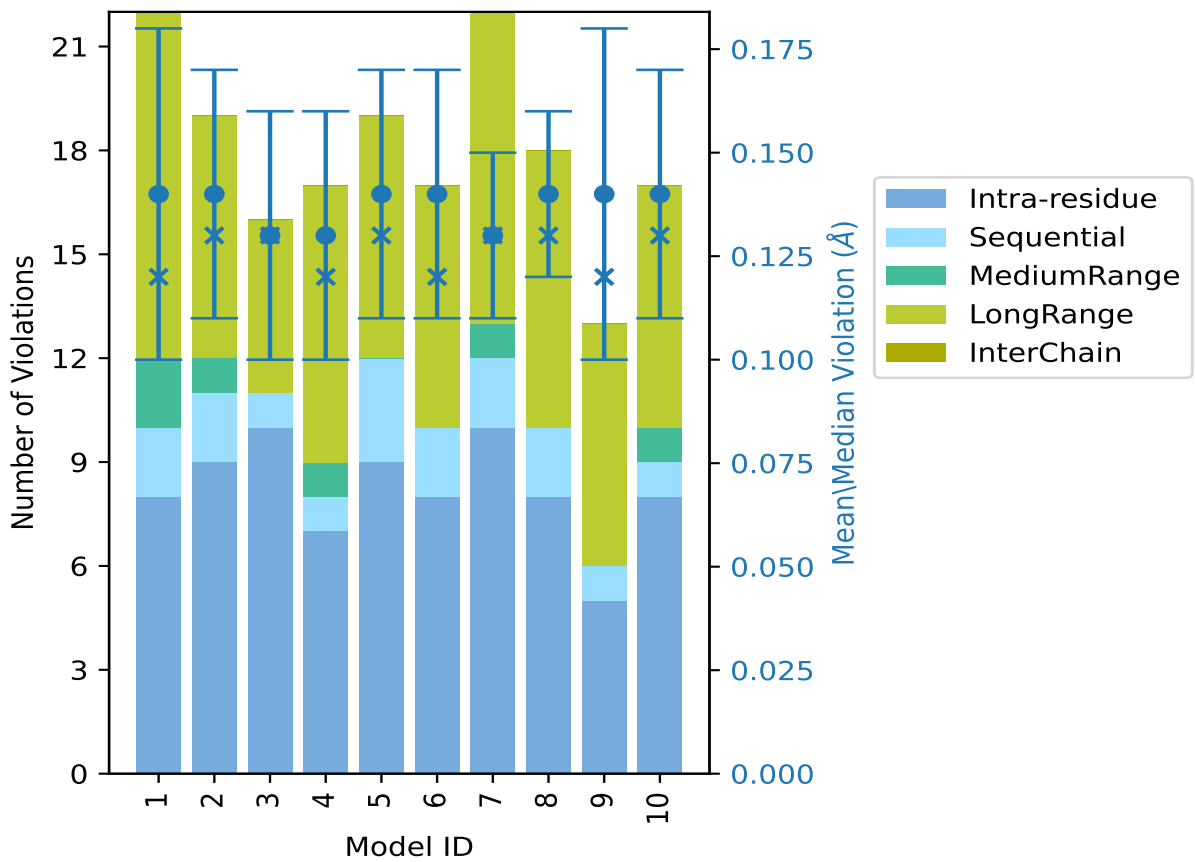
The following table provides the distance violation statistics for each model in the ensemble. Violations less than 0.1 Å are not included in the statistics.

Model ID	Number of violations						Mean (Å)	Max (Å)	SD[6] (Å)	Median (Å)
	IR[1]	SQ[2]	MR[3]	LR[4]	IC[5]	Total				
1	8	2	2	10	0	22	0.14	0.25	0.04	0.12
2	9	2	1	7	0	19	0.14	0.24	0.03	0.13
3	10	1	0	5	0	16	0.13	0.23	0.03	0.13
4	7	1	1	8	0	17	0.13	0.2	0.03	0.12
5	9	3	0	7	0	19	0.14	0.23	0.03	0.13
6	8	2	0	7	0	17	0.14	0.22	0.03	0.12
7	10	2	1	9	0	22	0.13	0.17	0.02	0.13
8	8	2	0	8	0	18	0.14	0.21	0.02	0.13
9	5	1	0	7	0	13	0.14	0.25	0.04	0.12
10	8	1	1	7	0	17	0.14	0.21	0.03	0.13

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

[5]Inter-chain restraints, [6]Standard 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 (i)

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, 1109(IR:532, SQ:206, MR:120, LR:251, 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]	%
4	1	2	7	0	14	1	10.0
2	1	2	0	0	5	2	20.0
1	1	0	4	0	6	3	30.0
2	1	0	2	0	5	4	40.0

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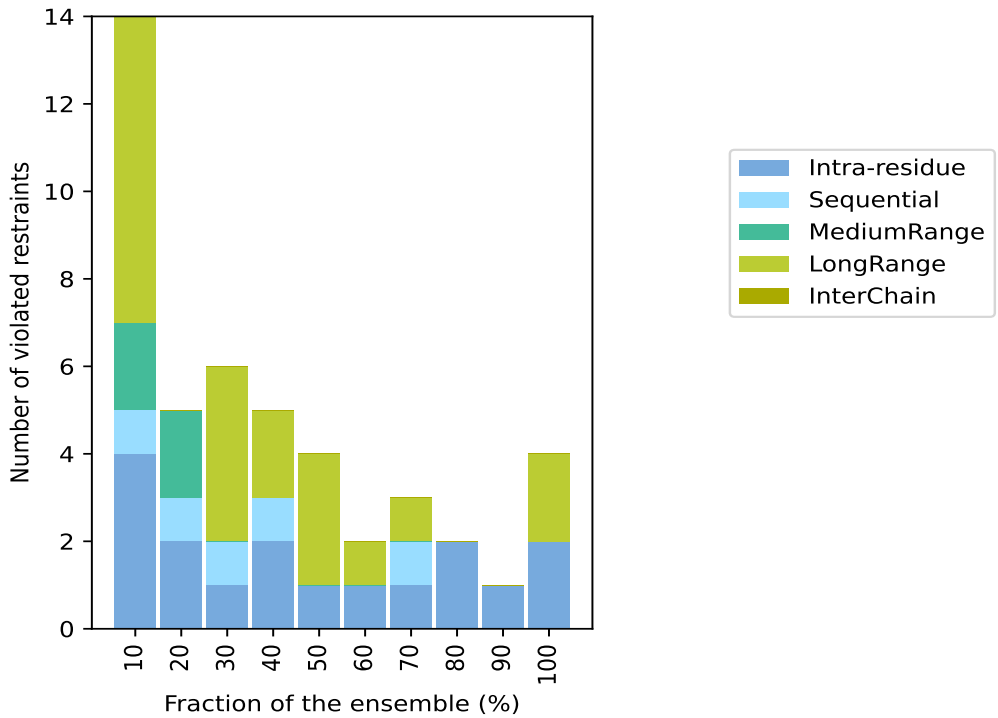
Number of violated restraints						Fraction of the ensemble	
IR[1]	SQ[2]	MR[3]	LR[4]	IC[5]	Total	Count[6]	%
1	0	0	3	0	4	5	50.0
1	0	0	1	0	2	6	60.0
1	1	0	1	0	3	7	70.0
2	0	0	0	0	2	8	80.0
1	0	0	0	0	1	9	90.0
2	0	0	2	0	4	10	100.0

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

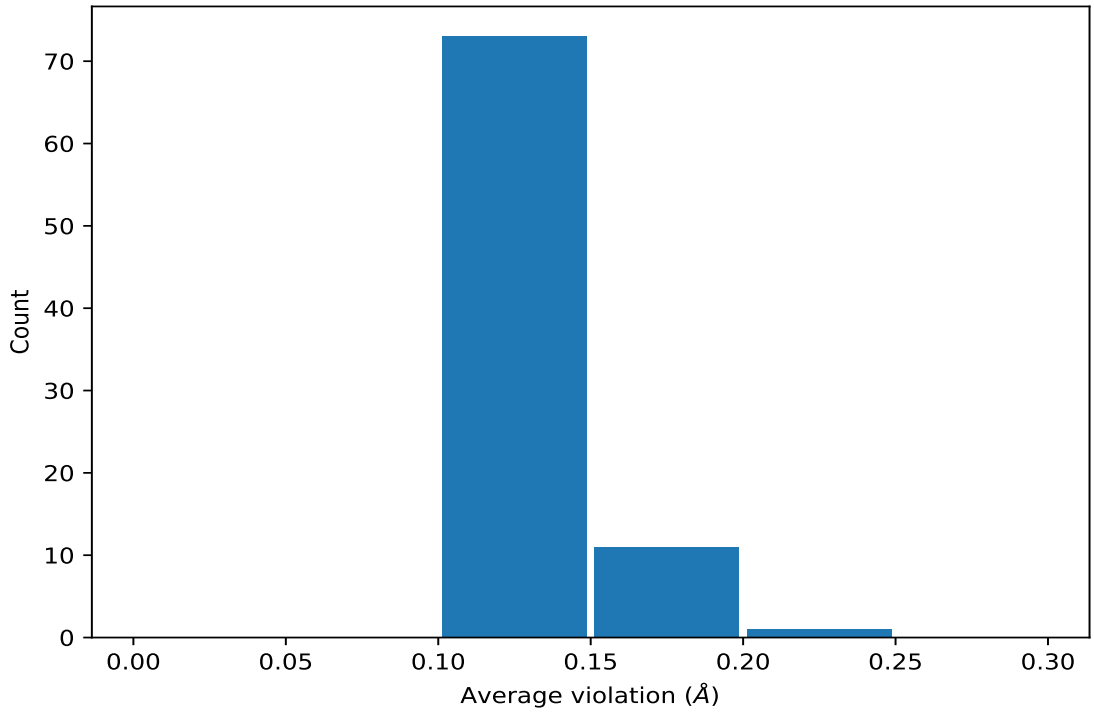
9.3.1 Bar graph : Distance violation statistics for the ensemble (i)



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 violations for the 10 worst performing restraints, sorted by number of violated models and the mean violation 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.

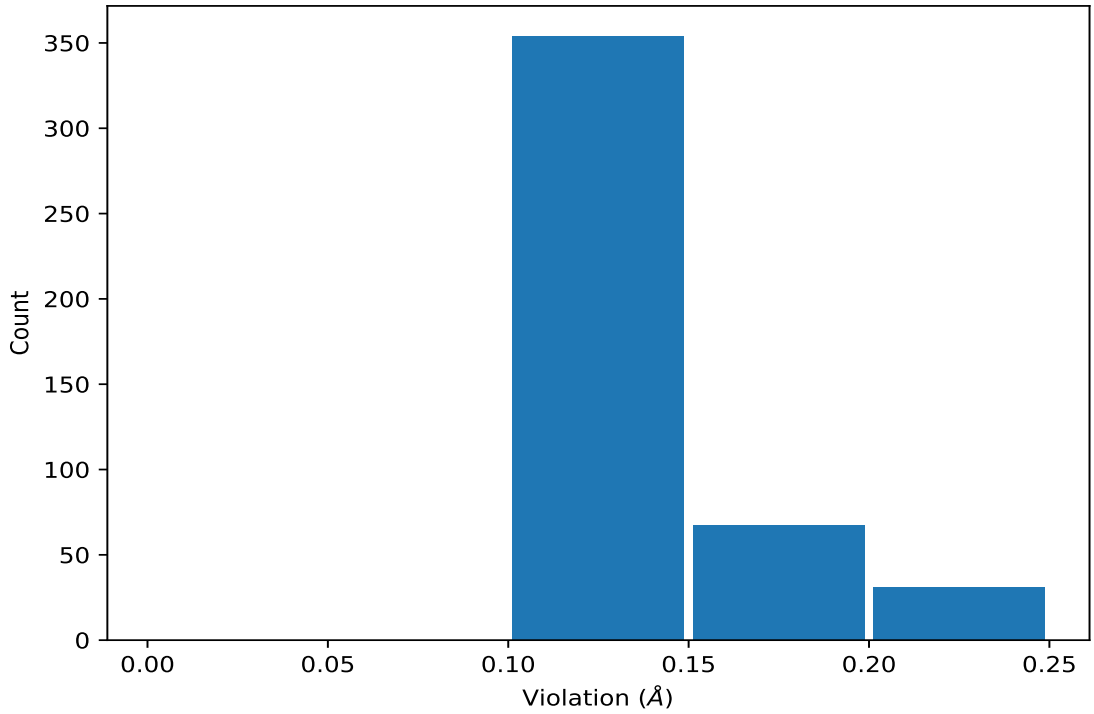
Key	Atom-1	Atom-2	Models[1]	Mean (Å)	SD[1] (Å)	Median (Å)
(1,892)	1:A:68:PHE:HE1	1:A:58:ALA:HB2	10	0.17	0.02	0.17
(1,892)	1:A:68:PHE:HE1	1:A:58:ALA:HB1	10	0.17	0.02	0.17
(1,892)	1:A:68:PHE:HE1	1:A:58:ALA:HB3	10	0.17	0.02	0.17
(1,892)	1:A:68:PHE:HE2	1:A:58:ALA:HB2	10	0.17	0.02	0.17
(1,892)	1:A:68:PHE:HE2	1:A:58:ALA:HB1	10	0.17	0.02	0.17
(1,892)	1:A:68:PHE:HE2	1:A:58:ALA:HB3	10	0.17	0.02	0.17
(1,769)	1:A:63:PHE:HE1	1:A:63:PHE:HZ	10	0.13	0.01	0.13
(1,927)	1:A:94:HIS:HE1	1:A:73:CYS:HB2	10	0.13	0.01	0.13
(1,933)	1:A:49:THR:HA	1:A:49:THR:HB	10	0.12	0.01	0.12
(1,916)	1:A:52:PRO:HA	1:A:52:PRO:HB2	9	0.11	0.0	0.11
(1,785)	1:A:78:SER:HB2	1:A:78:SER:HA	8	0.14	0.01	0.15
(1,264)	1:A:97:GLN:H	1:A:97:GLN:HB3	8	0.12	0.0	0.12
(1,737)	1:A:68:PHE:H	1:A:68:PHE:HD1	7	0.18	0.03	0.17
(1,1058)	1:A:60:LYS:H	1:A:59:ASP:HB2	7	0.18	0.04	0.17
(1,1058)	1:A:61:LEU:H	1:A:59:ASP:HB2	7	0.18	0.04	0.17
(1,1090)	1:A:67:CYS:H	1:A:79:LEU:HB2	7	0.14	0.02	0.14

[1]Number of violated models, [2]Standard deviation

9.5 All violated distance restraints (i)

9.5.1 Histogram : Distribution of distance violations (i)

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



9.5.2 Table : All distance violations (i)

The following table provides the 10 worst performing restraints, sorted by the violation 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)
(1,939)	1:A:60:LYS:HA	1:A:60:LYS:HB2	1	0.25
(1,1058)	1:A:60:LYS:H	1:A:59:ASP:HB2	9	0.25
(1,1058)	1:A:61:LEU:H	1:A:59:ASP:HB2	9	0.25
(1,719)	1:A:103:GLY:H	1:A:92:LYS:H	2	0.24
(1,737)	1:A:68:PHE:H	1:A:68:PHE:HD1	5	0.23
(1,437)	1:A:59:ASP:H	1:A:60:LYS:H	3	0.23
(1,737)	1:A:68:PHE:H	1:A:68:PHE:HD1	6	0.22
(1,1058)	1:A:60:LYS:H	1:A:59:ASP:HB2	1	0.22

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1058)	1:A:61:LEU:H	1:A:59:ASP:HB2	1	0.22
(1,892)	1:A:68:PHE:HE1	1:A:58:ALA:HB2	8	0.21
(1,892)	1:A:68:PHE:HE1	1:A:58:ALA:HB1	8	0.21
(1,892)	1:A:68:PHE:HE1	1:A:58:ALA:HB3	8	0.21
(1,892)	1:A:68:PHE:HE2	1:A:58:ALA:HB2	8	0.21
(1,892)	1:A:68:PHE:HE2	1:A:58:ALA:HB1	8	0.21
(1,892)	1:A:68:PHE:HE2	1:A:58:ALA:HB3	8	0.21
(1,737)	1:A:68:PHE:H	1:A:68:PHE:HD1	10	0.21
(1,437)	1:A:59:ASP:H	1:A:60:LYS:H	2	0.21

10 Dihedral-angle violation analysis (i)

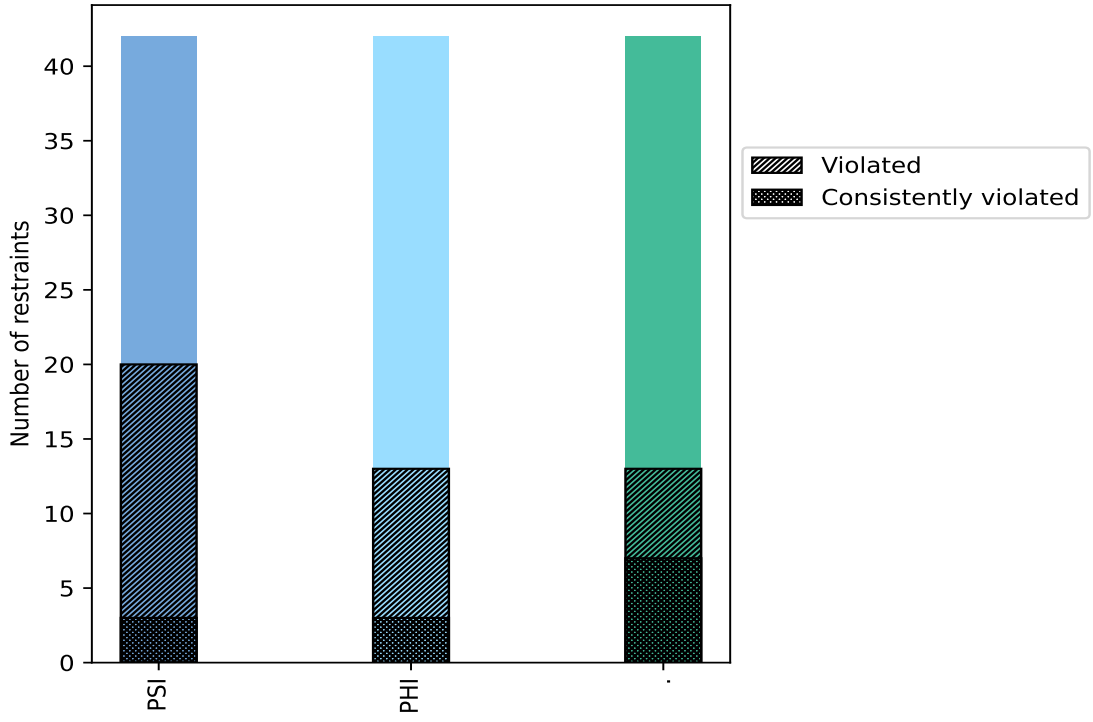
10.1 Summary of dihedral-angle violations (i)

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	42	33.3	20	47.6	15.9	3	7.1	2.4
PHI	42	33.3	13	31.0	10.3	3	7.1	2.4
.	42	33.3	13	31.0	10.3	7	16.7	5.6
Total	126	100.0	46	36.5	36.5	13	10.3	10.3

[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)



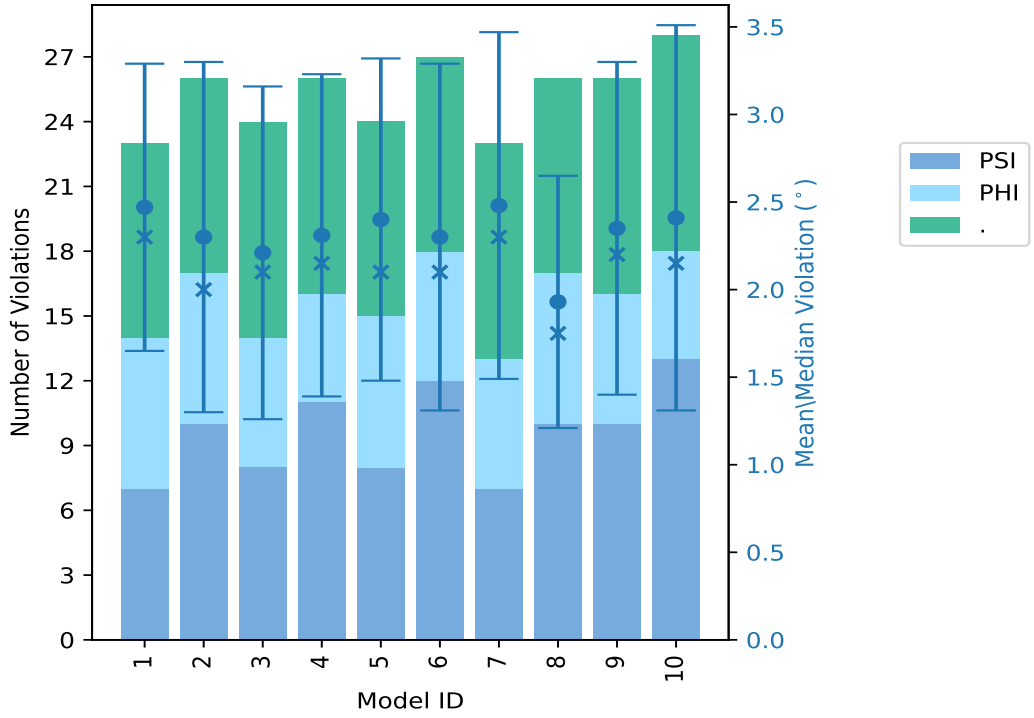
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

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	.	Total				
1	7	7	9	23	2.47	4.0	0.82	2.3
2	10	7	9	26	2.3	4.2	1.0	2.0
3	8	6	10	24	2.21	4.9	0.95	2.1
4	11	5	10	26	2.31	4.2	0.92	2.15
5	8	7	9	24	2.4	4.5	0.92	2.1
6	12	6	9	27	2.3	4.6	0.99	2.1
7	7	6	10	23	2.48	4.1	0.99	2.3
8	10	7	9	26	1.93	3.9	0.72	1.75
9	10	6	10	26	2.35	4.0	0.95	2.2
10	13	5	10	28	2.41	4.7	1.1	2.15

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

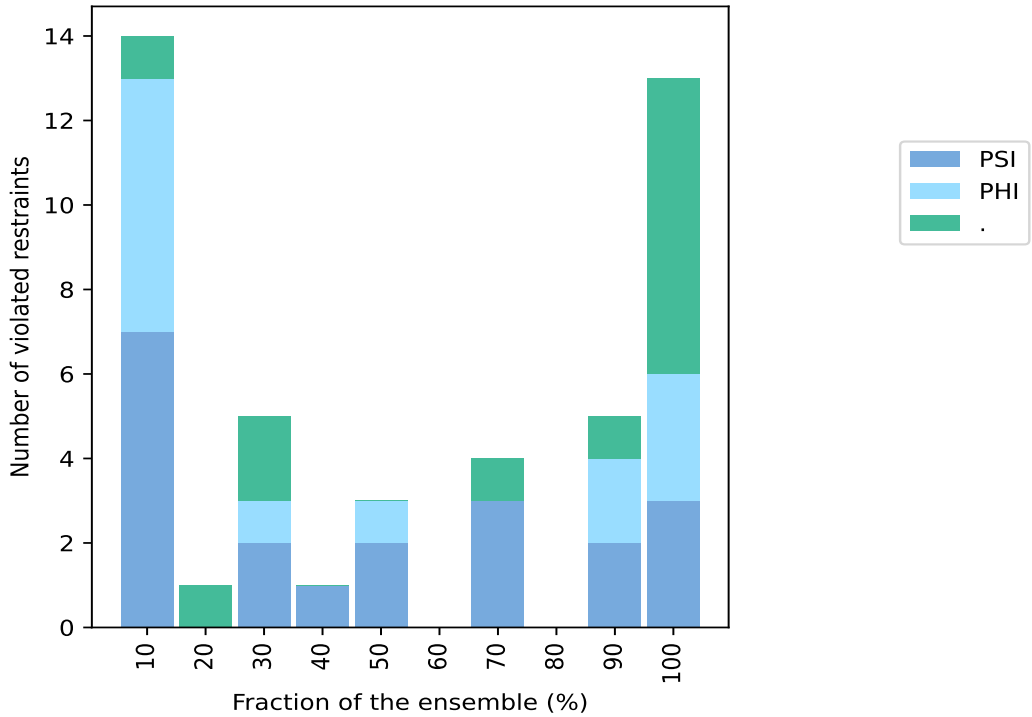
10.3 Dihedral-angle violation statistics for the ensemble (i)

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	.	Total	Count[1]	%
7	6	1	14	1	10.0
0	0	1	1	2	20.0
2	1	2	5	3	30.0
1	0	0	1	4	40.0
2	1	0	3	5	50.0
0	0	0	0	6	60.0
3	0	1	4	7	70.0
0	0	0	0	8	80.0
2	2	1	5	9	90.0
3	3	7	13	10	100.0

[1] Number of models with violations

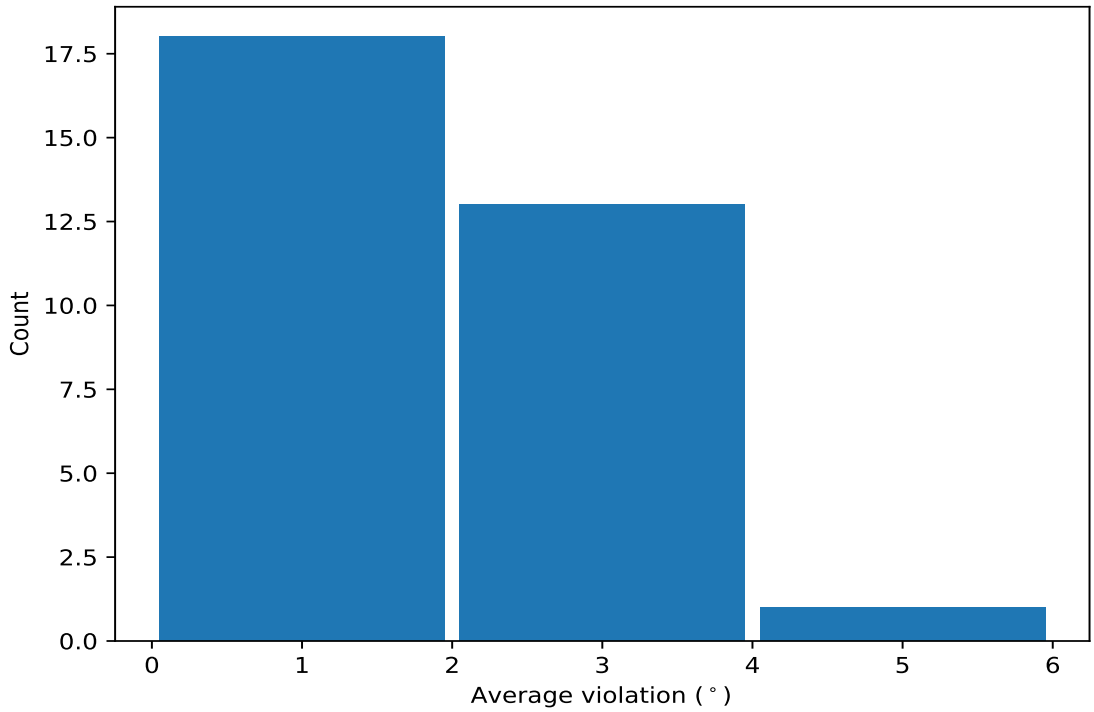
10.3.1 Bar graph : Dihedral-angle Violation statistics for the ensemble (i)



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 violations for the 10 worst performing restraints, sorted by number of violated models and the mean violation value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint.

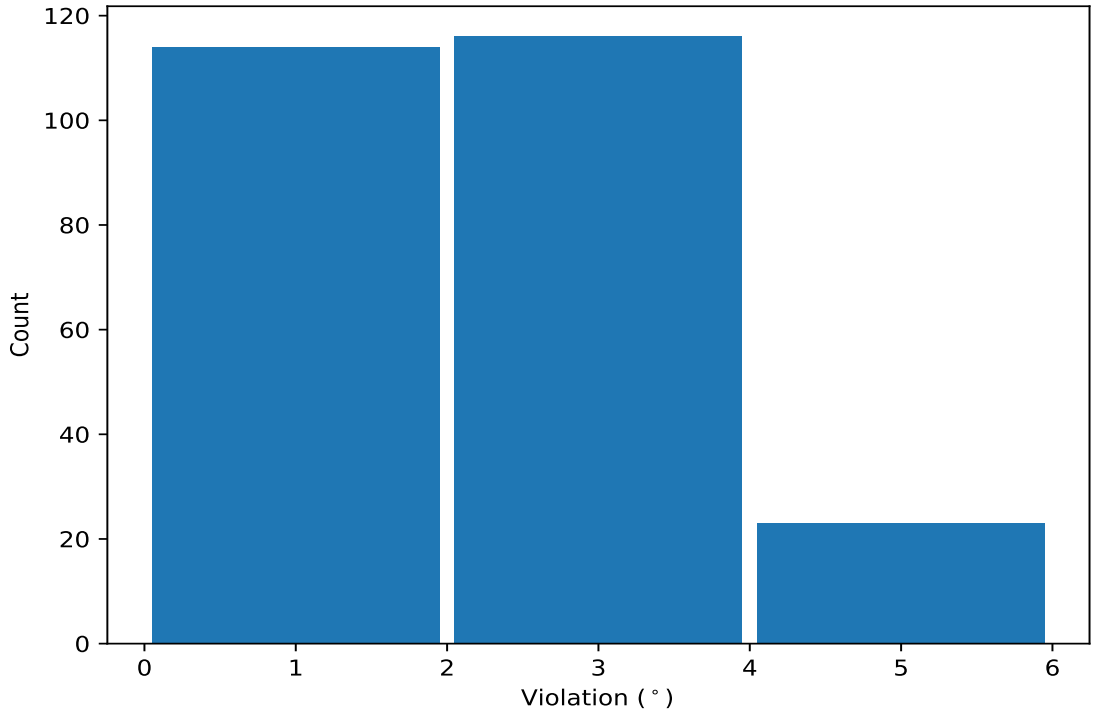
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models[1]	Mean	SD[2]	Median
(1,87)	1:A:46:CYS:N	1:A:46:CYS:CA	1:A:46:CYS:CB	1:A:46:CYS:HB2	10	4.03	0.09	4.0
(1,24)	1:A:58:ALA:N	1:A:58:ALA:CA	1:A:58:ALA:C	1:A:59:ASP:N	10	3.77	0.75	4.0
(1,48)	1:A:78:SER:N	1:A:78:SER:CA	1:A:78:SER:C	1:A:79:LEU:N	10	3.73	0.68	3.85
(1,81)	1:A:97:GLN:C	1:A:98:LEU:N	1:A:98:LEU:CA	1:A:98:LEU:C	10	3.4	0.86	3.45
(1,80)	1:A:97:GLN:N	1:A:97:GLN:CA	1:A:97:GLN:C	1:A:98:LEU:N	10	2.86	0.81	3.0
(1,107)	1:A:77:LEU:N	1:A:77:LEU:CA	1:A:77:LEU:CB	1:A:77:LEU:HB3	10	2.83	0.36	2.8
(1,33)	1:A:65:ASN:C	1:A:66:SER:N	1:A:66:SER:CA	1:A:66:SER:C	10	2.74	0.52	2.65
(1,116)	1:A:90:TYR:N	1:A:90:TYR:CA	1:A:90:TYR:CB	1:A:90:TYR:HB2	10	2.58	0.29	2.55
(1,95)	1:A:63:PHE:N	1:A:63:PHE:CA	1:A:63:PHE:CB	1:A:63:PHE:HB3	10	2.41	0.2	2.4
(1,100)	1:A:68:PHE:N	1:A:68:PHE:CA	1:A:68:PHE:CB	1:A:68:PHE:HB2	10	2.05	0.2	2.05

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

10.5 All violated dihedral-angle restraints (i)

10.5.1 Histogram : Distribution of violations (i)

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



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

The following table provides the list of violationsfor the 10 worst performing restraints, sorted by the violation 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 (°)
(1,48)	1:A:78:SER:N	1:A:78:SER:CA	1:A:78:SER:C	1:A:79:LEU:N	3	4.9
(1,81)	1:A:97:GLN:C	1:A:98:LEU:N	1:A:98:LEU:CA	1:A:98:LEU:C	10	4.7
(1,81)	1:A:97:GLN:C	1:A:98:LEU:N	1:A:98:LEU:CA	1:A:98:LEU:C	6	4.6
(1,24)	1:A:58:ALA:N	1:A:58:ALA:CA	1:A:58:ALA:C	1:A:59:ASP:N	10	4.6
(1,24)	1:A:58:ALA:N	1:A:58:ALA:CA	1:A:58:ALA:C	1:A:59:ASP:N	5	4.5
(1,87)	1:A:46:CYS:N	1:A:46:CYS:CA	1:A:46:CYS:CB	1:A:46:CYS:HB2	4	4.2
(1,48)	1:A:78:SER:N	1:A:78:SER:CA	1:A:78:SER:C	1:A:79:LEU:N	2	4.2
(1,24)	1:A:58:ALA:N	1:A:58:ALA:CA	1:A:58:ALA:C	1:A:59:ASP:N	2	4.2
(1,24)	1:A:58:ALA:N	1:A:58:ALA:CA	1:A:58:ALA:C	1:A:59:ASP:N	4	4.2
(1,87)	1:A:46:CYS:N	1:A:46:CYS:CA	1:A:46:CYS:CB	1:A:46:CYS:HB2	3	4.1