



# Full wwPDB NMR Structure Validation Report ⓘ

Dec 13, 2023 – 07:02 PM EST

PDB ID : 2JR7  
BMRB ID : 15377  
Title : Solution structure of human DESR1  
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Deposited on : 2007-06-21

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The types of validation reports are described at

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

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

Cyrange : Kirchner and Güntert (2011)  
NmrClust : Kelley et al. (1996)  
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) : 2.36

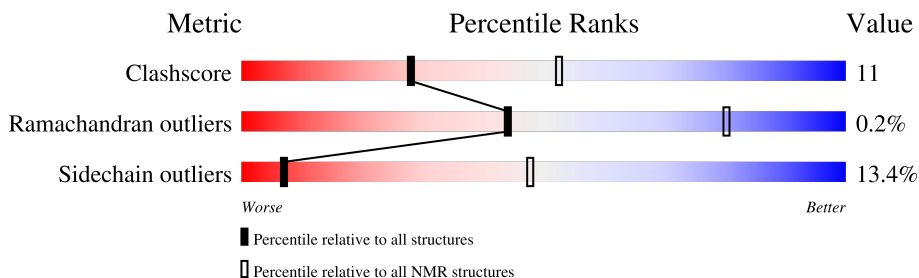
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 88%.

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  $\geq 3$ , 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	89	

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 9 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

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:4-A:63 (60)	0.32	9

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

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

The models can be grouped into 3 clusters and 1 single-model cluster was found.

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

### 3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1275 atoms, of which 609 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called DPH3 homolog.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	84	1274	419	609	98	143	5	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	66	SER	CYS	engineered mutation	UNP Q96FX2
A	82	LEU	-	expression tag	UNP Q96FX2
A	83	GLU	-	expression tag	UNP Q96FX2
A	84	HIS	-	expression tag	UNP Q96FX2
A	85	HIS	-	expression tag	UNP Q96FX2
A	86	HIS	-	expression tag	UNP Q96FX2
A	87	HIS	-	expression tag	UNP Q96FX2
A	88	HIS	-	expression tag	UNP Q96FX2
A	89	HIS	-	expression tag	UNP Q96FX2

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

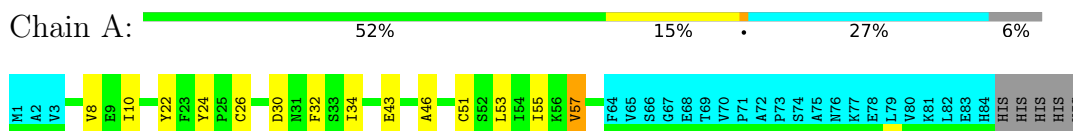
Mol	Chain	Residues	Atoms	
			Total	Zn
2	A	1	1	1

## 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: DPH3 homolog

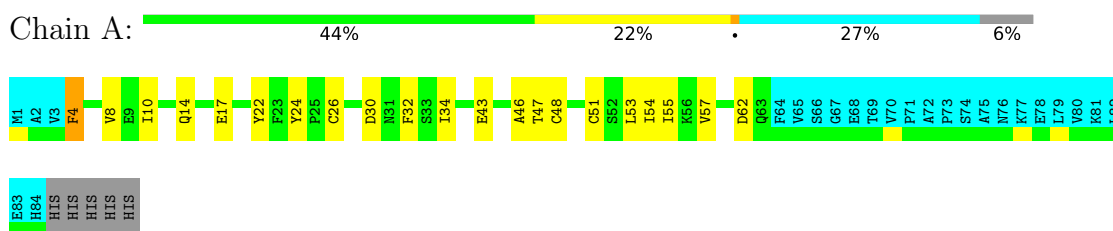


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

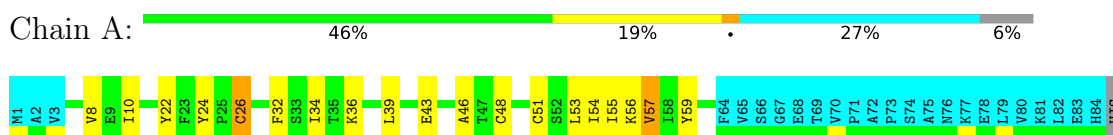
#### 4.2.1 Score per residue for model 1

- Molecule 1: DPH3 homolog



#### 4.2.2 Score per residue for model 2

- Molecule 1: DPH3 homolog



HIS  
HIS  
HIS  
HIS

### 4.2.3 Score per residue for model 3

- Molecule 1: DPH3 homolog

Chain A: 43% 22% 27% 6%

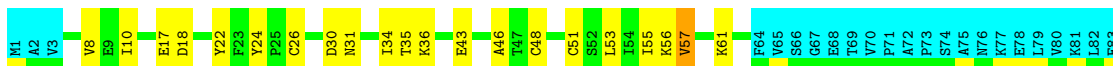


V80  
K81  
L82  
E83  
H84  
HIS  
HIS  
HIS  
HIS

### 4.2.4 Score per residue for model 4

- Molecule 1: DPH3 homolog

Chain A: 44% 22% 27% 6%



H84  
HIS  
HIS  
HIS  
HIS  
HIS

### 4.2.5 Score per residue for model 5

- Molecule 1: DPH3 homolog

Chain A: 44% 21% 27% 6%

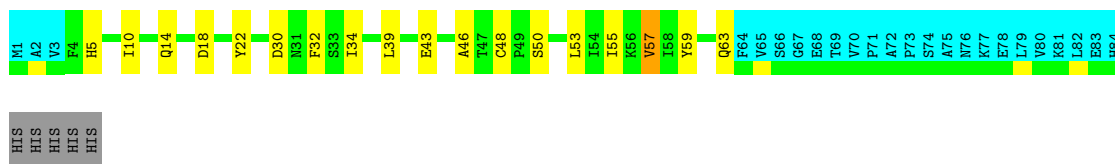


V80  
K81  
L82  
E83  
H84  
HIS  
HIS  
HIS  
HIS

### 4.2.6 Score per residue for model 6

- Molecule 1: DPH3 homolog

Chain A: 47% 19% 27% 6%



HIS  
HIS  
HIS  
HIS  
HIS

#### 4.2.7 Score per residue for model 7

- Molecule 1: DPH3 homolog

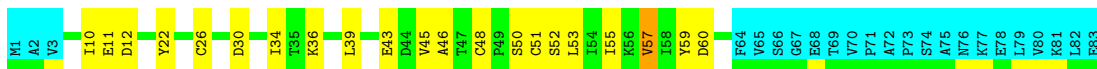
Chain A: 48% 16% 27% 6%



#### 4.2.8 Score per residue for model 8

- Molecule 1: DPH3 homolog

Chain A: 44% 22% 27% 6%



HIS  
HIS  
HIS  
HIS  
HIS

#### 4.2.9 Score per residue for model 9 (medoid)

- Molecule 1: DPH3 homolog

Chain A: 49% 16% 27% 6%

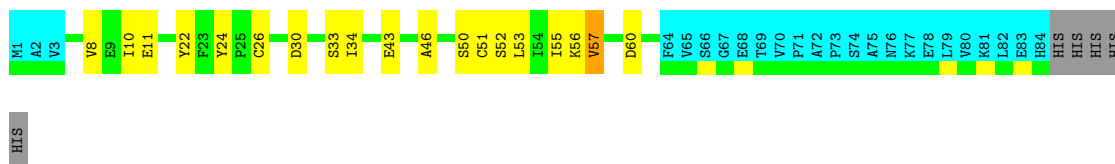


HIS

#### 4.2.10 Score per residue for model 10

- Molecule 1: DPH3 homolog

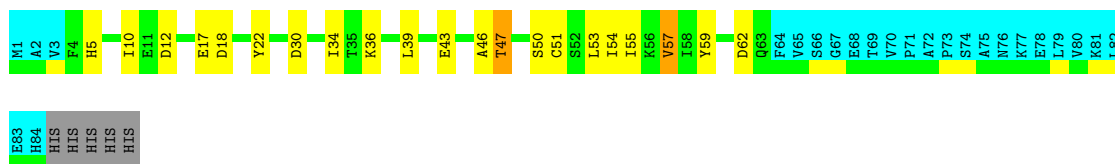
Chain A: 46% 20% 27% 6%



#### 4.2.11 Score per residue for model 11

- Molecule 1: DPH3 homolog

Chain A: 44% 21% 27% 6%



#### 4.2.12 Score per residue for model 12

- Molecule 1: DPH3 homolog

Chain A: 49% 16% 27% 6%



#### 4.2.13 Score per residue for model 13

- Molecule 1: DPH3 homolog

Chain A: 45% 20% 27% 6%

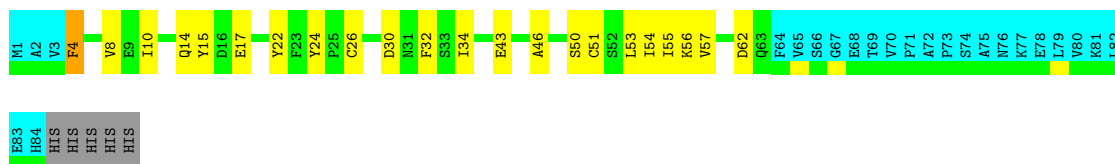


#### 4.2.14 Score per residue for model 14

- Molecule 1: DPH3 homolog

Chain A: 43% 24% 27% 6%





#### 4.2.15 Score per residue for model 15

- Molecule 1: DPH3 homolog

Chain A: 45% 22% 27% 6%



#### 4.2.16 Score per residue for model 16

- Molecule 1: DPH3 homolog

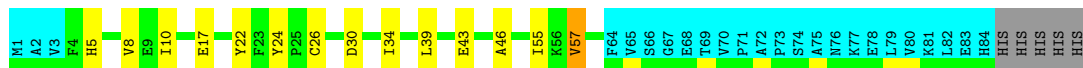
Chain A: 51% 16% 27% 6%



#### 4.2.17 Score per residue for model 17

- Molecule 1: DPH3 homolog

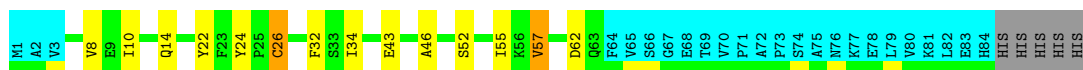
Chain A: 52% 15% 27% 6%



#### 4.2.18 Score per residue for model 18

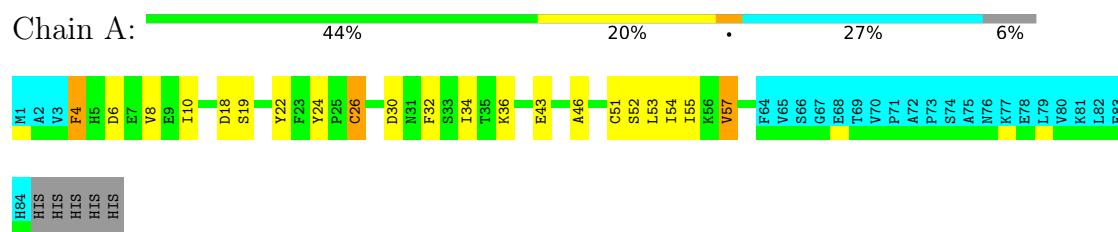
- Molecule 1: DPH3 homolog

Chain A: 52% 13% 27% 6%



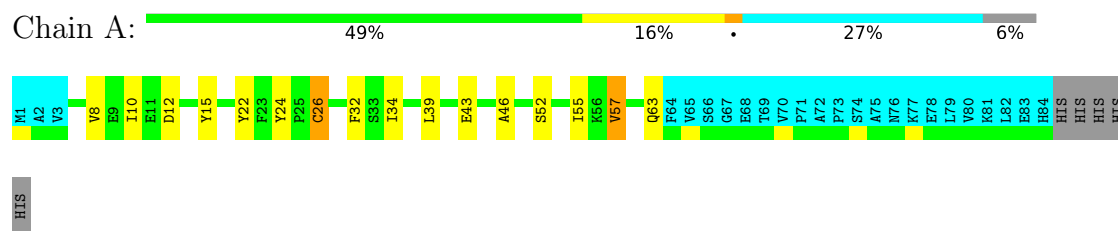
### 4.2.19 Score per residue for model 19

- Molecule 1: DPH3 homolog



### 4.2.20 Score per residue for model 20

- Molecule 1: DPH3 homolog



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *torsion angle dynamics*.

Of the 200 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
CNS	structure solution	1.1
CNS	refinement	1.1

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	928
Number of shifts mapped to atoms	925
Number of unparsed shifts	0
Number of shifts with mapping errors	3
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	88%

## 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	487	424	423	10±2
All	All	9760	8480	8460	193

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 11.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:34:ILE:HD11	1:A:43:GLU:O	0.72	1.84	5	20
1:A:10:ILE:HD11	1:A:22:TYR:CD2	0.71	2.21	2	17
1:A:34:ILE:HD13	1:A:57:VAL:HG21	0.70	1.63	4	17
1:A:8:VAL:HG11	1:A:24:TYR:CZ	0.68	2.24	17	5
1:A:10:ILE:HD11	1:A:22:TYR:CG	0.67	2.25	5	16
1:A:46:ALA:HB3	1:A:55:ILE:CG1	0.64	2.21	14	19
1:A:51:CYS:SG	1:A:53:LEU:HD13	0.64	2.33	11	3
1:A:8:VAL:HG11	1:A:24:TYR:CE1	0.63	2.29	1	11
1:A:32:PHE:HB3	1:A:46:ALA:HB1	0.58	1.75	14	5
1:A:4:PHE:CD1	1:A:54:ILE:HD12	0.58	2.33	3	4
1:A:39:LEU:HD22	1:A:59:TYR:CE2	0.57	2.35	2	2
1:A:46:ALA:HB3	1:A:55:ILE:HD11	0.56	1.76	7	12

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:46:ALA:HB3	1:A:55:ILE:CD1	0.52	2.35	3	11
1:A:51:CYS:SG	1:A:53:LEU:HD12	0.50	2.46	4	9
1:A:46:ALA:HB3	1:A:55:ILE:HG13	0.49	1.83	5	9
1:A:5:HIS:CD2	1:A:53:LEU:HD21	0.49	2.43	5	1
1:A:10:ILE:HD11	1:A:22:TYR:CD1	0.48	2.43	20	4
1:A:8:VAL:HG21	1:A:24:TYR:CZ	0.45	2.46	4	1
1:A:10:ILE:HD13	1:A:39:LEU:HD13	0.45	1.88	8	4
1:A:5:HIS:CG	1:A:53:LEU:HD21	0.44	2.48	6	2
1:A:39:LEU:HD22	1:A:59:TYR:CE1	0.43	2.48	3	2
1:A:26:CYS:SG	1:A:32:PHE:CE2	0.43	3.12	13	9
1:A:15:TYR:HB2	1:A:22:TYR:CE2	0.43	2.49	14	1
1:A:47:THR:HG23	1:A:53:LEU:O	0.42	2.14	11	1
1:A:39:LEU:HD22	1:A:59:TYR:CD1	0.42	2.49	3	2
1:A:15:TYR:CD1	1:A:22:TYR:CZ	0.41	3.08	14	3
1:A:8:VAL:HG21	1:A:24:TYR:CE2	0.41	2.50	18	1
1:A:10:ILE:HD13	1:A:39:LEU:CD1	0.41	2.45	8	1
1:A:51:CYS:HB2	1:A:53:LEU:HD13	0.40	1.92	3	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	60/89 (67%)	52±1 (86±2%)	8±1 (14±2%)	0±0 (0±1%)	44 80
All	All	1200/1780 (67%)	1030 (86%)	167 (14%)	3 (0%)	44 80

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	31	ASN	2
1	A	50	SER	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	57/82 (70%)	49±2 (87±3%)	8±2 (13±3%)	<b>7</b> 48
All	All	1140/1640 (70%)	987 (87%)	153 (13%)	<b>7</b> 48

All 29 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	57	VAL	20
1	A	26	CYS	18
1	A	30	ASP	15
1	A	50	SER	9
1	A	36	LYS	8
1	A	52	SER	8
1	A	14	GLN	7
1	A	17	GLU	7
1	A	18	ASP	7
1	A	48	CYS	6
1	A	62	ASP	6
1	A	56	LYS	5
1	A	47	THR	4
1	A	11	GLU	4
1	A	4	PHE	3
1	A	63	GLN	3
1	A	12	ASP	3
1	A	60	ASP	3
1	A	37	GLU	2
1	A	61	LYS	2
1	A	10	ILE	2
1	A	45	VAL	2
1	A	19	SER	2
1	A	5	HIS	2
1	A	35	THR	1
1	A	33	SER	1
1	A	7	GLU	1
1	A	16	ASP	1
1	A	6	ASP	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](#)

Of 1 ligands modelled in this entry, 1 is 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

The completeness of assignment taking into account all chemical shift lists is 88% for the well-defined parts and 86% for the entire structure.

### 7.1 Chemical shift list 1

File name: working\_cs.cif

Chemical shift list name: *assigned\_chem\_shift\_list\_1*

#### 7.1.1 Bookkeeping

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

Total number of shifts	928
Number of shifts mapped to atoms	925
Number of unparsed shifts	0
Number of shifts with mapping errors	3
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. All 3 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	89	HIS	H	8.286	0.011	.
1	A	89	HIS	C	174.836	0.000	.
1	A	89	HIS	N	119.932	0.018	.

#### 7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	81	-0.22 $\pm$ 0.11	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	78	-0.03 $\pm$ 0.15	None needed (< 0.5 ppm)
$^{13}\text{C}'$	81	-0.01 $\pm$ 0.16	None needed (< 0.5 ppm)
$^{15}\text{N}$	76	-0.84 $\pm$ 0.31	Should be applied



### 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 88%, i.e. 674 atoms were assigned a chemical shift out of a possible 766. 0 out of 5 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	<b>Total</b>	<b><sup>1</sup>H</b>	<b><sup>13</sup>C</b>	<b><sup>15</sup>N</b>
Backbone	296/296 (100%)	119/119 (100%)	120/120 (100%)	57/57 (100%)
Sidechain	348/386 (90%)	237/245 (97%)	111/134 (83%)	0/7 (0%)
Aromatic	30/84 (36%)	30/40 (75%)	0/42 (0%)	0/2 (0%)
Overall	674/766 (88%)	386/404 (96%)	231/296 (78%)	57/66 (86%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 86%, i.e. 925 atoms were assigned a chemical shift out of a possible 1079. 0 out of 11 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	<b>Total</b>	<b><sup>1</sup>H</b>	<b><sup>13</sup>C</b>	<b><sup>15</sup>N</b>
Backbone	395/413 (96%)	159/166 (96%)	161/168 (96%)	75/79 (95%)
Sidechain	496/564 (88%)	341/363 (94%)	155/191 (81%)	0/10 (0%)
Aromatic	34/102 (33%)	34/49 (69%)	0/49 (0%)	0/4 (0%)
Overall	925/1079 (86%)	534/578 (92%)	316/408 (77%)	75/93 (81%)

### 7.1.4 Statistically unusual chemical shifts [i](#)

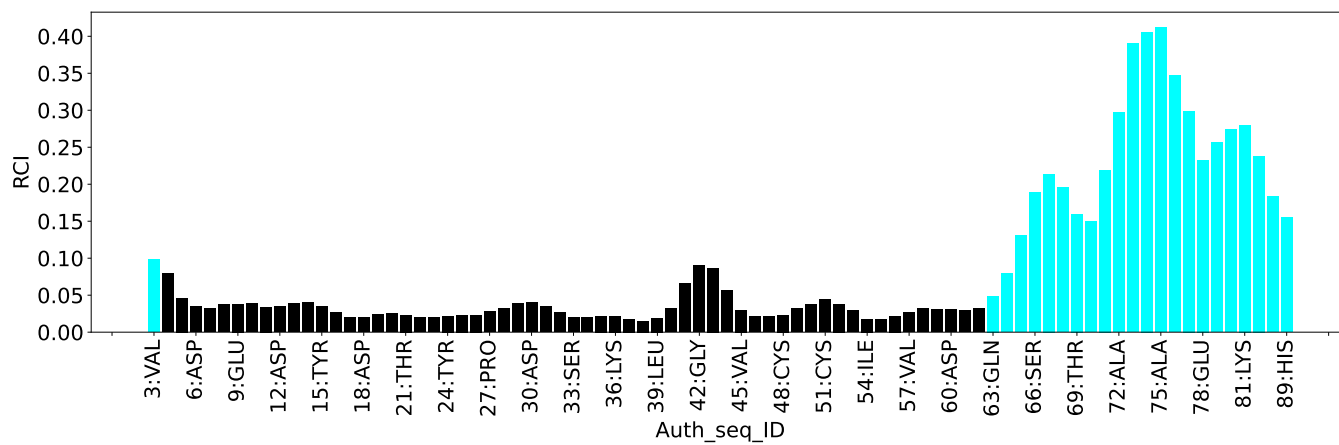
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	56	LYS	HB3	0.13	0.46 – 3.04	-6.3

### 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	1567
Intra-residue ( $ i-j =0$ )	279
Sequential ( $ i-j =1$ )	453
Medium range ( $ i-j >1$ and $ i-j <5$ )	239
Long range ( $ i-j \geq 5$ )	566
Inter-chain	0
Hydrogen bond restraints	20
Disulfide bond restraints	6
Total dihedral-angle restraints	60
Number of unmapped restraints	0
Number of restraints per residue	18.1
Number of long range restraints per residue <sup>1</sup>	6.6

<sup>1</sup>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

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)	1.1	0.17
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)	0.1	1.13
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

## 9 Distance violation analysis

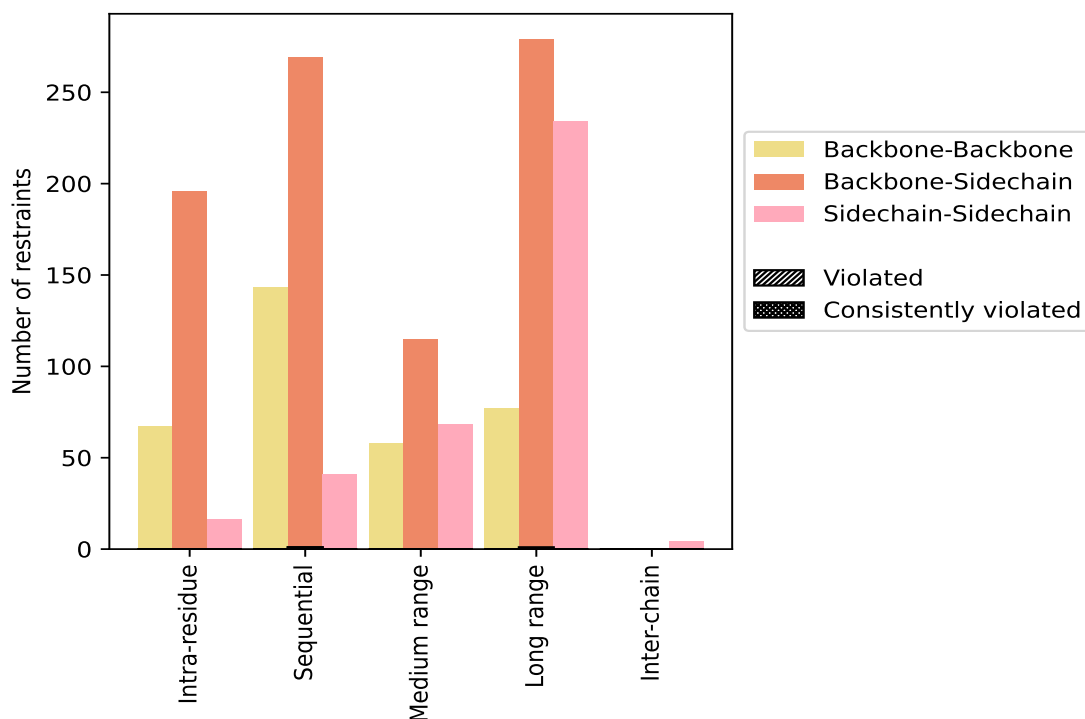
### 9.1 Summary of distance violations

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

Restrains type	Count	% <sup>1</sup>	Violated <sup>3</sup>			Consistently Violated <sup>4</sup>		
			Count	% <sup>2</sup>	% <sup>1</sup>	Count	% <sup>2</sup>	% <sup>1</sup>
<b>Intra-residue (<math> i-j =0</math>)</b>	<b>279</b>	<b>17.8</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	67	4.3	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	196	12.5	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	16	1.0	0	0.0	0.0	0	0.0	0.0
<b>Sequential (<math> i-j =1</math>)</b>	<b>453</b>	<b>28.9</b>	<b>1</b>	<b>0.2</b>	<b>0.1</b>	<b>1</b>	<b>0.2</b>	<b>0.1</b>
Backbone-Backbone	143	9.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	269	17.2	1	0.4	0.1	1	0.4	0.1
Sidechain-Sidechain	41	2.6	0	0.0	0.0	0	0.0	0.0
<b>Medium range (<math> i-j &gt;1</math> &amp; <math> i-j &lt;5</math>)</b>	<b>239</b>	<b>15.3</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	58	3.7	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	115	7.3	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	66	4.2	0	0.0	0.0	0	0.0	0.0
<b>Long range (<math> i-j \geq 5</math>)</b>	<b>566</b>	<b>36.1</b>	<b>1</b>	<b>0.2</b>	<b>0.1</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
Backbone-Backbone	77	4.9	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	259	16.5	1	0.4	0.1	0	0.0	0.0
Sidechain-Sidechain	230	14.7	0	0.0	0.0	0	0.0	0.0
<b>Inter-chain</b>	<b>0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
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
<b>Hydrogen bond</b>	<b>20</b>	<b>1.3</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
<b>Disulfide bond</b>	<b>6</b>	<b>0.4</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>	<b>0</b>	<b>0.0</b>	<b>0.0</b>
<b>Total</b>	<b>1567</b>	<b>100.0</b>	<b>2</b>	<b>0.1</b>	<b>0.1</b>	<b>1</b>	<b>0.1</b>	<b>0.1</b>
Backbone-Backbone	345	22.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	859	54.8	2	0.2	0.1	1	0.1	0.1
Sidechain-Sidechain	363	23.2	0	0.0	0.0	0	0.0	0.0

<sup>1</sup> percentage calculated with respect to the total number of distance restraints, <sup>2</sup> percentage calculated with respect to the number of restraints in a particular restraint category, <sup>3</sup> violated in at least one model, <sup>4</sup> 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 <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total				
1	0	1	0	0	0	1	0.15	0.15	0.0	0.15
2	0	1	0	0	0	1	0.16	0.16	0.0	0.16
3	0	1	0	0	0	1	0.16	0.16	0.0	0.16
4	0	1	0	0	0	1	0.15	0.15	0.0	0.15
5	0	1	0	0	0	1	0.16	0.16	0.0	0.16
6	0	1	0	0	0	1	0.17	0.17	0.0	0.17
7	0	1	0	0	0	1	0.16	0.16	0.0	0.16
8	0	1	0	0	0	1	0.16	0.16	0.0	0.16
9	0	1	0	0	0	1	0.15	0.15	0.0	0.15
10	0	1	0	0	0	1	0.15	0.15	0.0	0.15
11	0	1	0	0	0	1	0.16	0.16	0.0	0.16

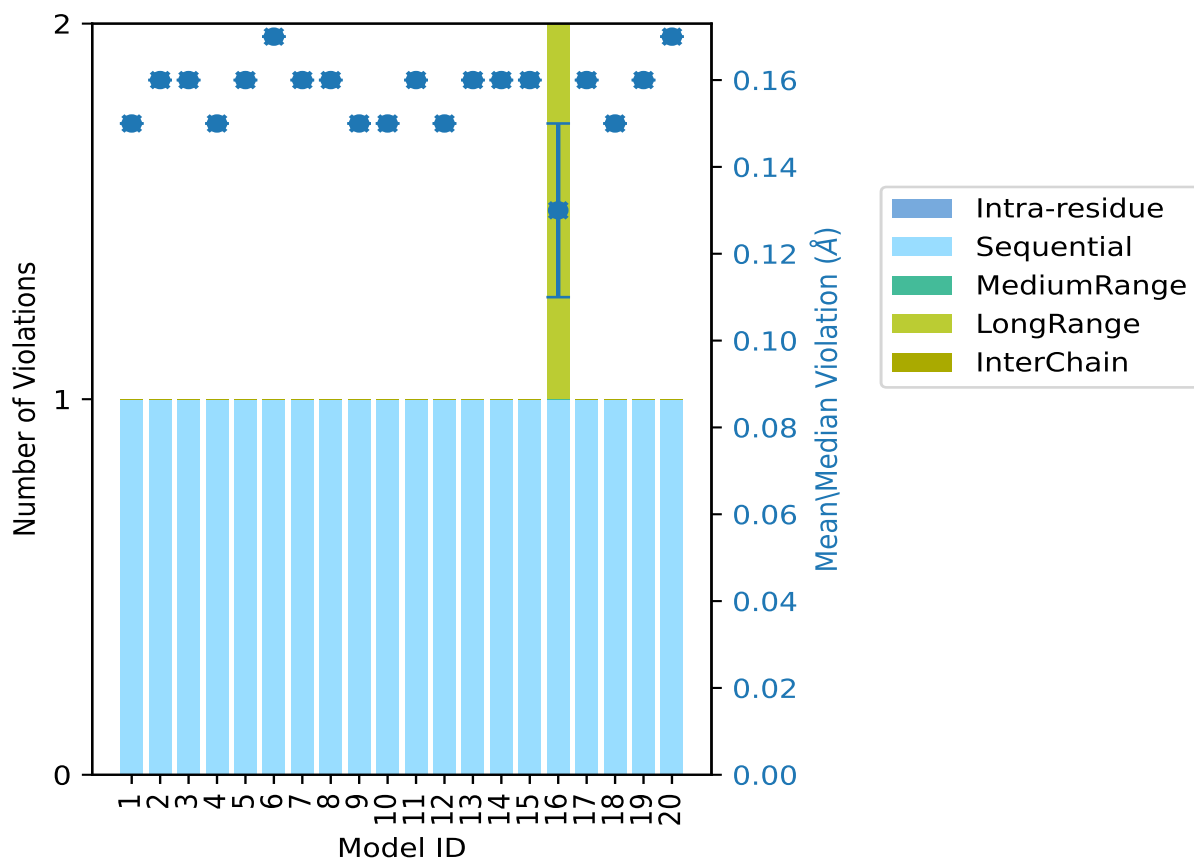
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD <sup>6</sup> (Å)	Median (Å)
	IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total				
12	0	1	0	0	0	1	0.15	0.15	0.0	0.15
13	0	1	0	0	0	1	0.16	0.16	0.0	0.16
14	0	1	0	0	0	1	0.16	0.16	0.0	0.16
15	0	1	0	0	0	1	0.16	0.16	0.0	0.16
16	0	1	0	1	0	2	0.13	0.15	0.02	0.13
17	0	1	0	0	0	1	0.16	0.16	0.0	0.16
18	0	1	0	0	0	1	0.15	0.15	0.0	0.15
19	0	1	0	0	0	1	0.16	0.16	0.0	0.16
20	0	1	0	0	0	1	0.17	0.17	0.0	0.17

<sup>1</sup>Intra-residue restraints, <sup>2</sup>Sequential restraints, <sup>3</sup>Medium range restraints, <sup>4</sup>Long range restraints, <sup>5</sup>Inter-chain restraints, <sup>6</sup>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

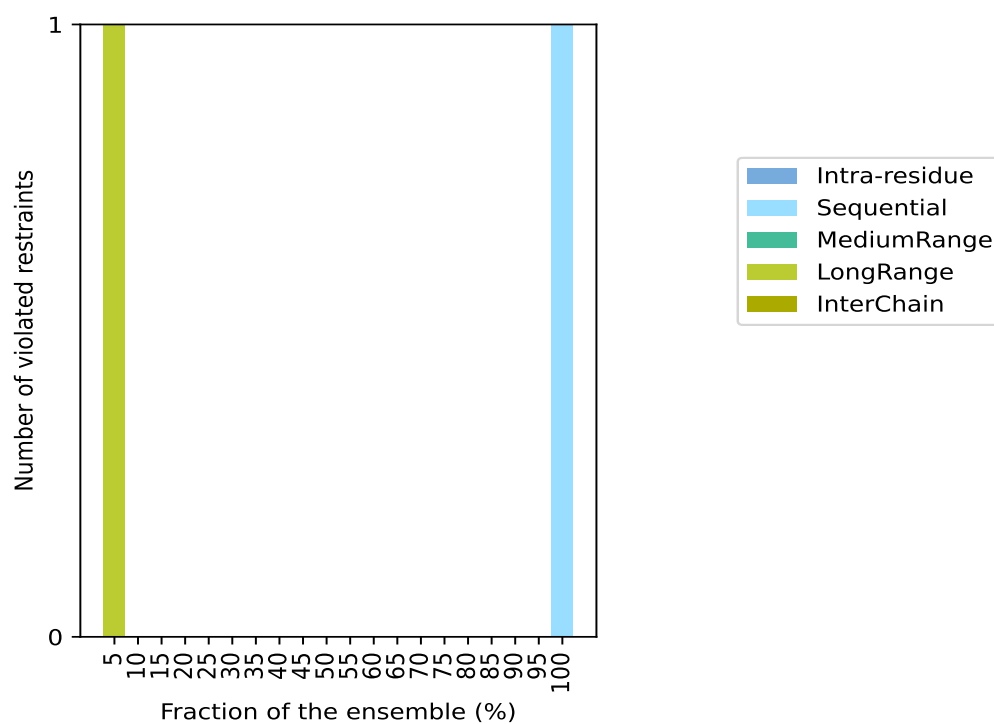
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 1535(IR:279, SQ:452, MR:239, LR:565, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR <sup>1</sup>	SQ <sup>2</sup>	MR <sup>3</sup>	LR <sup>4</sup>	IC <sup>5</sup>	Total	Count <sup>6</sup>	%
0	0	0	1	0	1	1	5.0
0	0	0	0	0	0	2	10.0
0	0	0	0	0	0	3	15.0
0	0	0	0	0	0	4	20.0
0	0	0	0	0	0	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
0	0	0	0	0	0	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	1	0	0	0	1	20	100.0

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



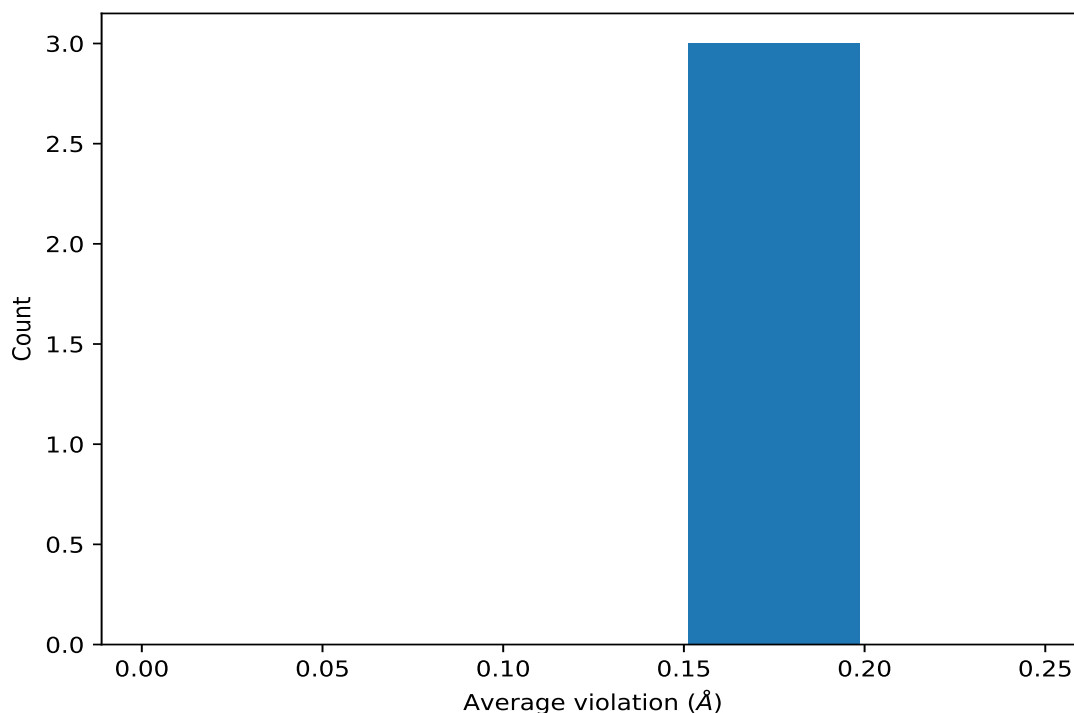
### 9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



## 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.

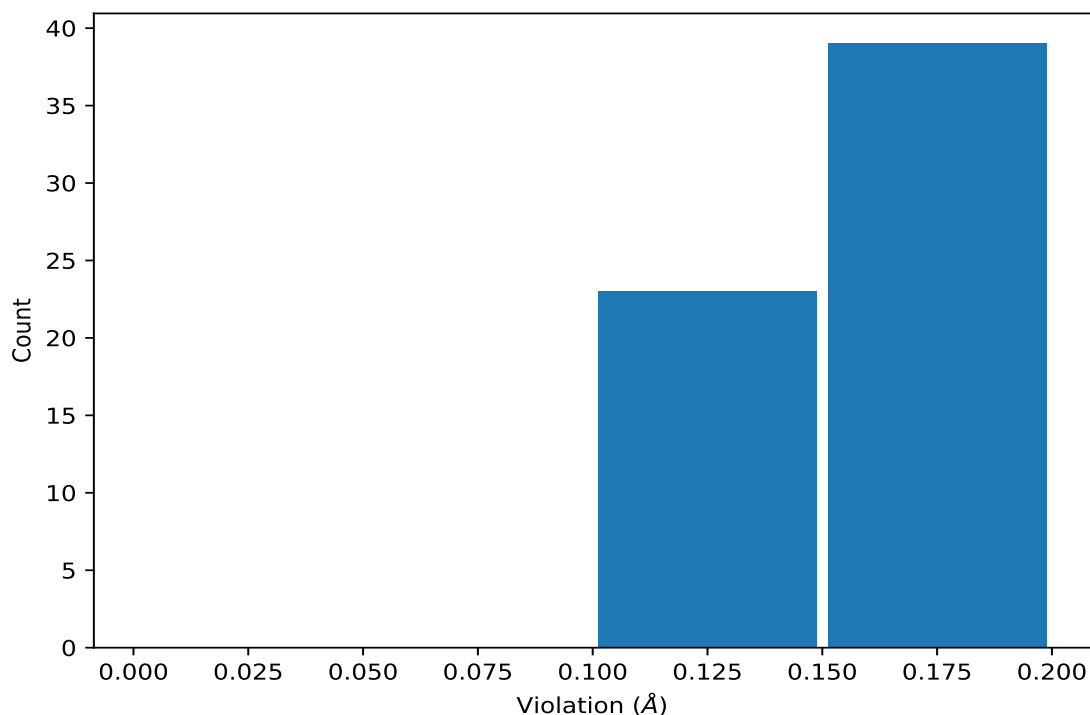
Key	Atom-1	Atom-2	Models <sup>1</sup>	Mean (Å)	SD <sup>1</sup> (Å)	Median (Å)
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	20	0.16	0.01	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	20	0.16	0.01	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	20	0.16	0.01	0.16

<sup>1</sup>Number of violated models, <sup>2</sup>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 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)
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	6	0.17
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	6	0.17
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	6	0.17
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	20	0.17
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	20	0.17
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	20	0.17
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	2	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	2	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	2	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	3	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	3	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	3	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	5	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	5	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	5	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	7	0.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	7	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	7	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	8	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	8	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	8	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	11	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	11	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	11	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	13	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	13	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	13	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	14	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	14	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	14	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	15	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	15	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	15	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	17	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	17	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	17	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	19	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	19	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	19	0.16
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	1	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	1	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	1	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	4	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	4	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	4	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	9	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	9	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	9	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	10	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	10	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	10	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	12	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	12	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	12	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	16	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	16	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	16	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG11	18	0.15

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<b>Key</b>	<b>Atom-1</b>	<b>Atom-2</b>	<b>Model ID</b>	<b>Violation (Å)</b>
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG12	18	0.15
(1,1007)	1:56:A:LYS:HA	1:57:A:VAL:HG13	18	0.15
(1,1462)	1:44:A:ASP:HA	1:59:A:TYR:HE1	16	0.11
(1,1462)	1:44:A:ASP:HA	1:59:A:TYR:HE2	16	0.11

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

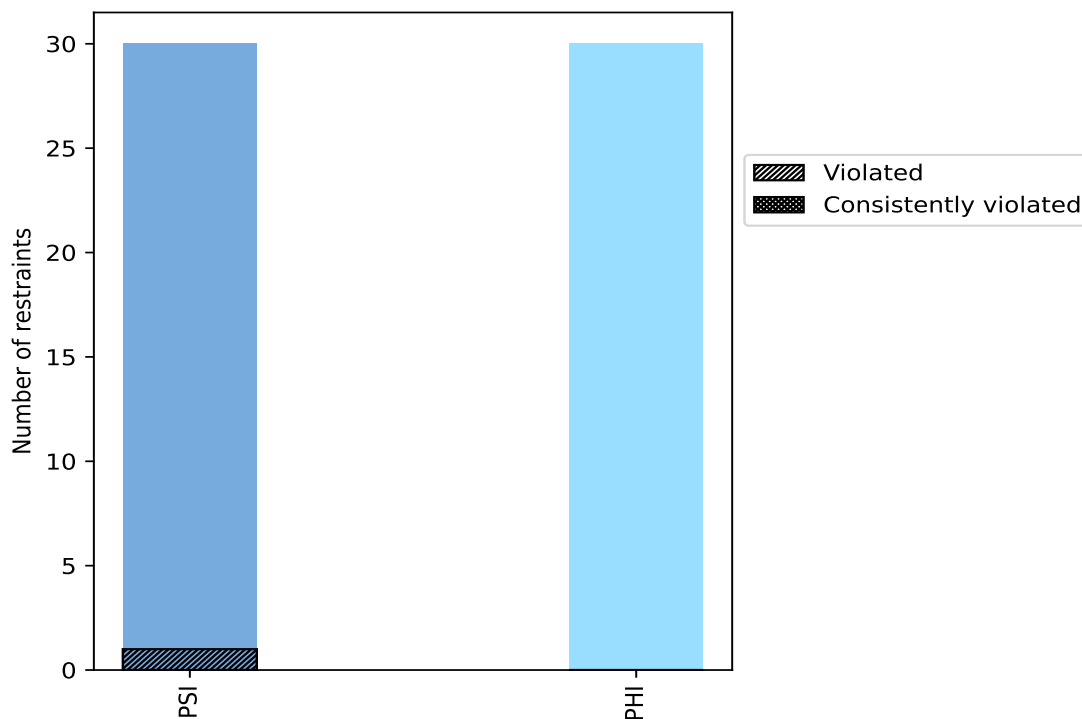
### 10.1 Summary of dihedral-angle violations [i](#)

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	% <sup>1</sup>	Violated <sup>3</sup>			Consistently Violated <sup>4</sup>		
			Count	% <sup>2</sup>	% <sup>1</sup>	Count	% <sup>2</sup>	% <sup>1</sup>
PSI	30	50.0	1	3.3	1.7	0	0.0	0.0
PHI	30	50.0	0	0.0	0.0	0	0.0	0.0
Total	60	100.0	1	1.7	1.7	0	0.0	0.0

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

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



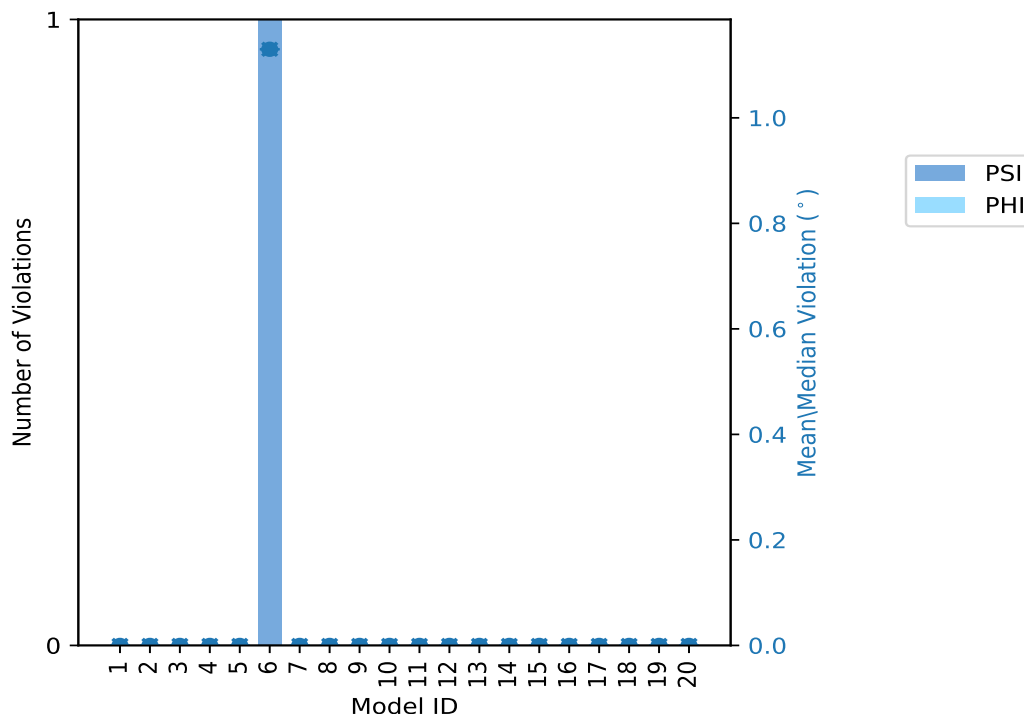
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

## 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	0	0	0	0.0	0.0	0.0	0.0
2	0	0	0	0.0	0.0	0.0	0.0
3	0	0	0	0.0	0.0	0.0	0.0
4	0	0	0	0.0	0.0	0.0	0.0
5	0	0	0	0.0	0.0	0.0	0.0
6	1	0	1	1.13	1.13	0.0	1.13
7	0	0	0	0.0	0.0	0.0	0.0
8	0	0	0	0.0	0.0	0.0	0.0
9	0	0	0	0.0	0.0	0.0	0.0
10	0	0	0	0.0	0.0	0.0	0.0
11	0	0	0	0.0	0.0	0.0	0.0
12	0	0	0	0.0	0.0	0.0	0.0
13	0	0	0	0.0	0.0	0.0	0.0
14	0	0	0	0.0	0.0	0.0	0.0
15	0	0	0	0.0	0.0	0.0	0.0
16	0	0	0	0.0	0.0	0.0	0.0
17	0	0	0	0.0	0.0	0.0	0.0
18	0	0	0	0.0	0.0	0.0	0.0
19	0	0	0	0.0	0.0	0.0	0.0
20	0	0	0	0.0	0.0	0.0	0.0

### 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 <sup>1</sup>	%
1	0	1	1	5.0
0	0	0	2	10.0
0	0	0	3	15.0
0	0	0	4	20.0
0	0	0	5	25.0
0	0	0	6	30.0
0	0	0	7	35.0
0	0	0	8	40.0
0	0	0	9	45.0
0	0	0	10	50.0
0	0	0	11	55.0

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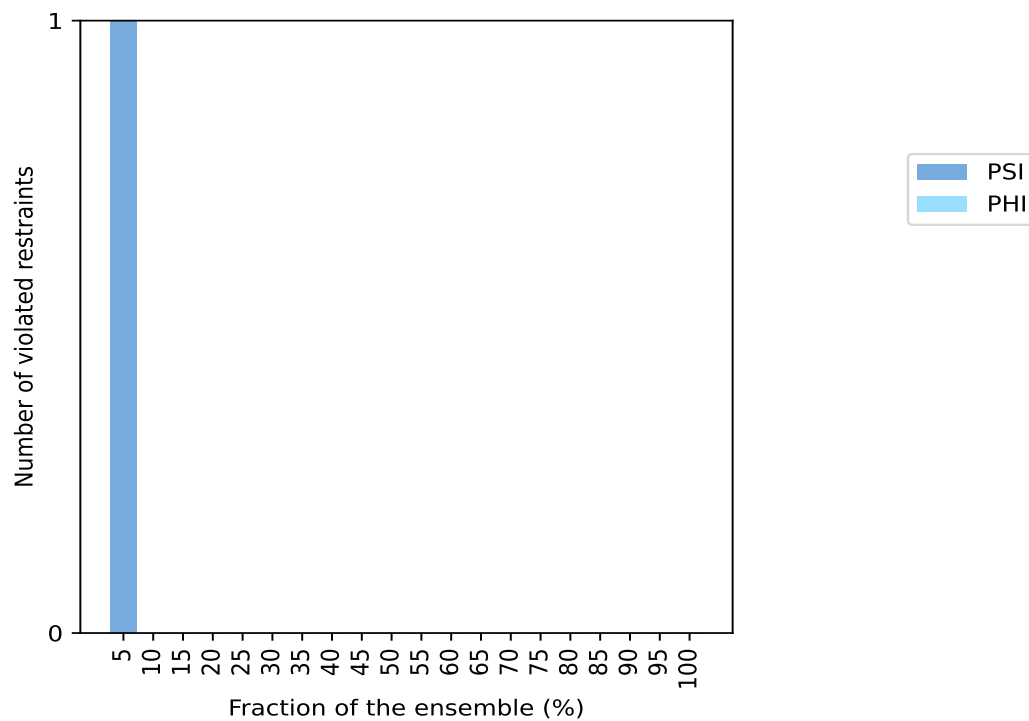


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Number of violated restraints			Fraction of the ensemble	
PSI	PHI	Total	Count <sup>1</sup>	%
0	0	0	12	60.0
0	0	0	13	65.0
0	0	0	14	70.0
0	0	0	15	75.0
0	0	0	16	80.0
0	0	0	17	85.0
0	0	0	18	90.0
0	0	0	19	95.0
0	0	0	20	100.0

<sup>1</sup> 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](#)

No violations found

## 10.5 All violated dihedral-angle restraints

### 10.5.1 Histogram : Distribution of violations

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

Data insufficient to plot histogram

### 10.5.2 Table: All violated dihedral-angle restraints

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 (°)
(1,46)	1:54:A:ILE:N	1:54:A:ILE:CA	1:54:A:ILE:C	1:55:A:ILE:N	6	1.13