Full wwPDB NMR Structure Validation Report (i)

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PDB ID	:	2E2F
BMRB ID	:	6729
Title	:	Solution structure of DSP
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This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (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 (i)) were used in the production of this report:

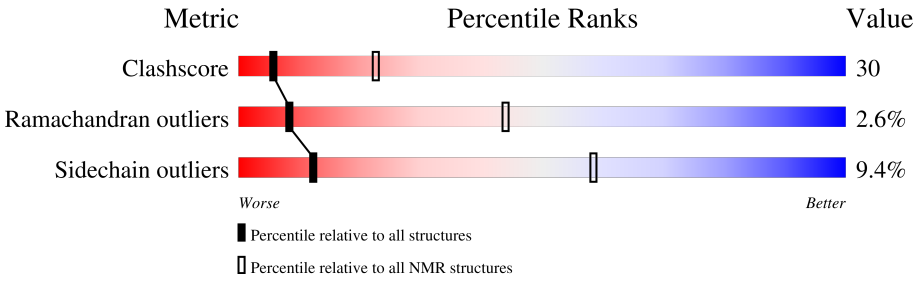
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
BMRB Restraints Analysis	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal 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 46%.

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	41	56% 39%

2 Ensemble composition and analysis (i)

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 592 atoms, of which 284 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Diapausin.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	41	592	184	284	62	55	7	0

4 Residue-property plots (i)

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: Diapausin

Chain A:

5 Refinement protocol and experimental data overview

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

Of the 50 calculated structures, 1 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
X-PLOR	refinement	3.1F
NMRPipe	structure solution	2.2
PIPP	structure solution	4.3.2
NMRView	structure solution	5.0.3
CYANA	structure solution	2.1
X-PLOR	structure solution	3.1F

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	236
Number of shifts mapped to atoms	236
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	46%

6 Model quality

6.1 Standard geometry

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

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

Mol	Chain	Chirality	Planarity
1	A	0	5
All	All	0	5

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

All planar outliers are listed below.

Mol	Chain	Res	Type	Group
1	A	3	ARG	Sidechain
1	A	13	ARG	Sidechain
1	A	18	ARG	Sidechain
1	A	23	ARG	Sidechain
1	A	27	ARG	Sidechain

6.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	308	284	284	18
All	All	308	284	284	18

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 30.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:4:ILE:HD11	1:A:27:ARG:HD3	0.68	1.65
1:A:7:CYS:HA	1:A:10:VAL:HG12	0.64	1.70
1:A:10:VAL:HG11	1:A:21:CYS:HB2	0.61	1.70
1:A:10:VAL:HG23	1:A:13:ARG:HD2	0.55	1.77
1:A:4:ILE:HB	1:A:7:CYS:SG	0.54	2.42
1:A:4:ILE:HG21	1:A:21:CYS:SG	0.53	2.43
1:A:4:ILE:HD12	1:A:25:HIS:CB	0.53	2.34
1:A:4:ILE:HB	1:A:7:CYS:HB2	0.53	1.80
1:A:4:ILE:HD12	1:A:25:HIS:HB3	0.49	1.83
1:A:10:VAL:HG23	1:A:13:ARG:CD	0.48	2.38
1:A:4:ILE:HB	1:A:7:CYS:CB	0.48	2.38
1:A:11:CYS:HB3	1:A:12:PRO:HD3	0.47	1.87
1:A:15:VAL:N	1:A:16:PRO:HD2	0.44	2.26
1:A:11:CYS:N	1:A:12:PRO:HD2	0.44	2.27
1:A:11:CYS:N	1:A:12:PRO:CD	0.42	2.82
1:A:7:CYS:HA	1:A:10:VAL:CG1	0.42	2.43
1:A:7:CYS:HB3	1:A:38:MET:HB3	0.41	1.93
1:A:4:ILE:HD11	1:A:27:ARG:CD	0.41	2.43

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	39/41 (95%)	31 (79%)	7 (18%)	1 (3%)	8 44
All	All	39/41 (95%)	31 (79%)	7 (18%)	1 (3%)	8 44

All 1 Ramachandran outliers are listed below.

Mol	Chain	Res	Type
1	A	14	ILE

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	32/32 (100%)	29 (91%)	3 (9%)	12 58
All	All	32/32 (100%)	29 (91%)	3 (9%)	12 58

All 3 residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	20	GLU
1	A	27	ARG
1	A	34	SER

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

The completeness of assignment taking into account all chemical shift lists is 46% for the well-defined parts and 46% 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	236
Number of shifts mapped to atoms	236
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

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

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 46%, i.e. 236 atoms were assigned a chemical shift out of a possible 515. 0 out of 3 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	78/205 (38%)	78/85 (92%)	0/82 (0%)	0/38 (0%)
Sidechain	146/269 (54%)	146/174 (84%)	0/78 (0%)	0/17 (0%)
Aromatic	12/41 (29%)	12/20 (60%)	0/19 (0%)	0/2 (0%)
Overall	236/515 (46%)	236/279 (85%)	0/179 (0%)	0/57 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 46%, i.e. 236 atoms were assigned a chemical shift out of a possible 515. 0 out of 3 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	78/205 (38%)	78/85 (92%)	0/82 (0%)	0/38 (0%)
Sidechain	146/269 (54%)	146/174 (84%)	0/78 (0%)	0/17 (0%)
Aromatic	12/41 (29%)	12/20 (60%)	0/19 (0%)	0/2 (0%)
Overall	236/515 (46%)	236/279 (85%)	0/179 (0%)	0/57 (0%)

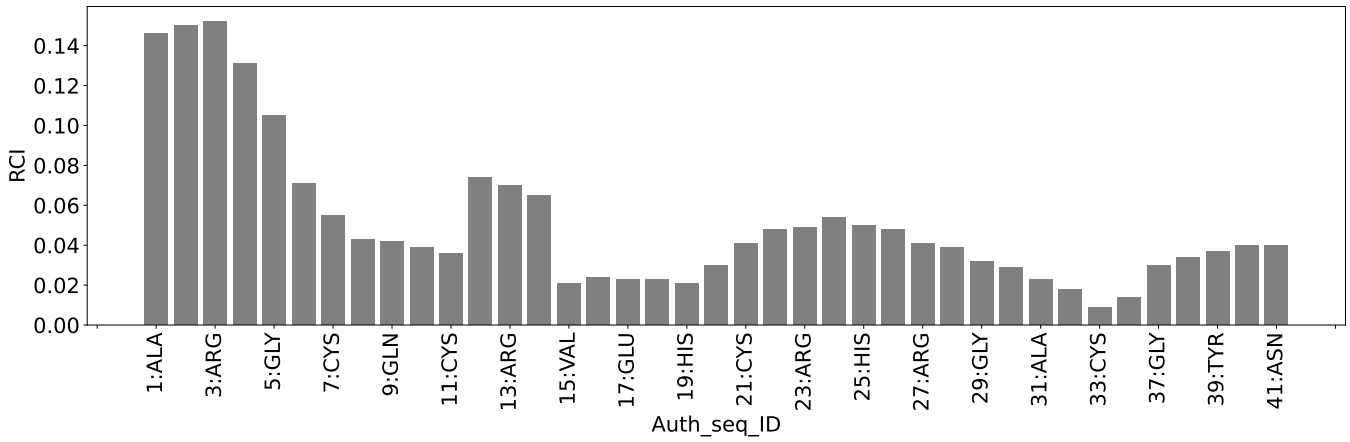
7.1.4 Statistically unusual chemical shifts (i)

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:



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	414
Intra-residue (|i-j|=0)	195
Sequential (|i-j|=1)	95
Medium range (|i-j|>1 and |i-j|<5)	55
Long range (|i-j|≥5)	63
Inter-chain	0
Hydrogen bond restraints	6
Disulfide bond restraints	0
Total dihedral-angle restraints	24
Number of unmapped restraints	0
Number of restraints per residue	10.7
Number of long range restraints per residue[1]	1.7

[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

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)	1.0	0.16
0.2-0.5 (Medium)	3.0	0.4
>0.5 (Large)	1.0	5.32

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)	5.0	8.1
10.0-20.0 (Medium)	4.0	16.8
>20.0 (Large)	3.0	44.6

9 Distance violation analysis (i)

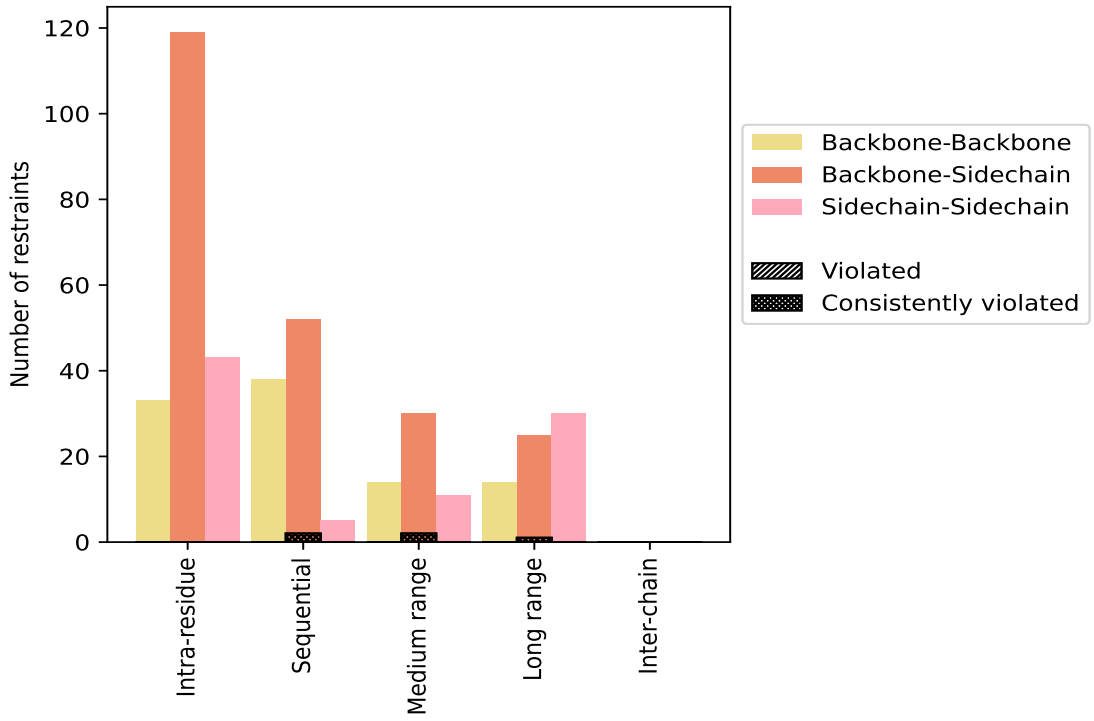
9.1 Summary of distance violations (i)

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)	195	47.1	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	33	8.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	119	28.7	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	43	10.4	0	0.0	0.0	0	0.0	0.0
Sequential (|i-j|=1)	95	22.9	2	2.1	0.5	2	2.1	0.5
Backbone-Backbone	38	9.2	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	52	12.6	2	3.8	0.5	2	3.8	0.5
Sidechain-Sidechain	5	1.2	0	0.0	0.0	0	0.0	0.0
Medium range (|i-j|>1 & |i-j|<5)	55	13.3	2	3.6	0.5	2	3.6	0.5
Backbone-Backbone	14	3.4	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	30	7.2	2	6.7	0.5	2	6.7	0.5
Sidechain-Sidechain	11	2.7	0	0.0	0.0	0	0.0	0.0
Long range (|i-j|≥5)	63	15.2	1	1.6	0.2	1	1.6	0.2
Backbone-Backbone	8	1.9	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	25	6.0	1	4.0	0.2	1	4.0	0.2
Sidechain-Sidechain	30	7.2	0	0.0	0.0	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	6	1.4	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	414	100.0	5	1.2	1.2	5	1.2	1.2
Backbone-Backbone	99	23.9	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	226	54.6	5	2.2	1.2	5	2.2	1.2
Sidechain-Sidechain	89	21.5	0	0.0	0.0	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



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

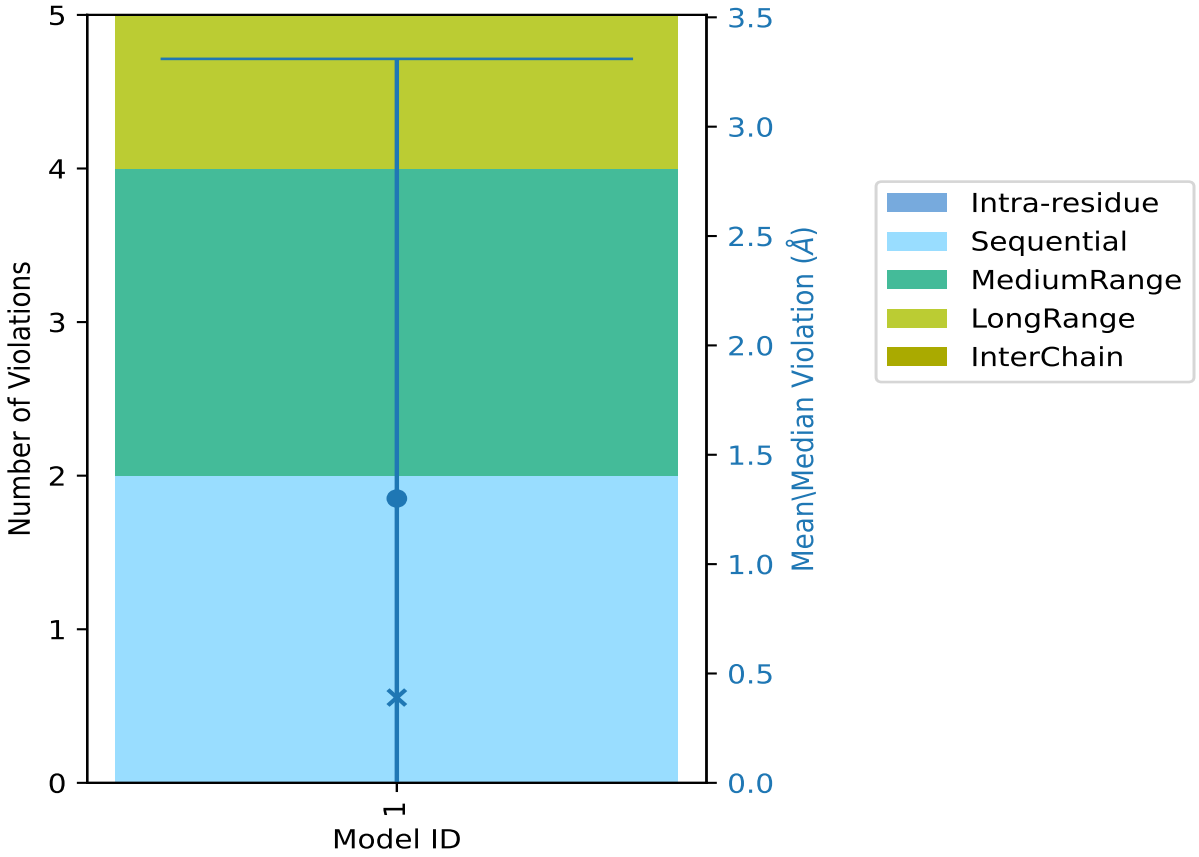
9.2 Distance violation statistics for each model

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	2	2	1	0	5	1.3	5.32	2.01	0.39

[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



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

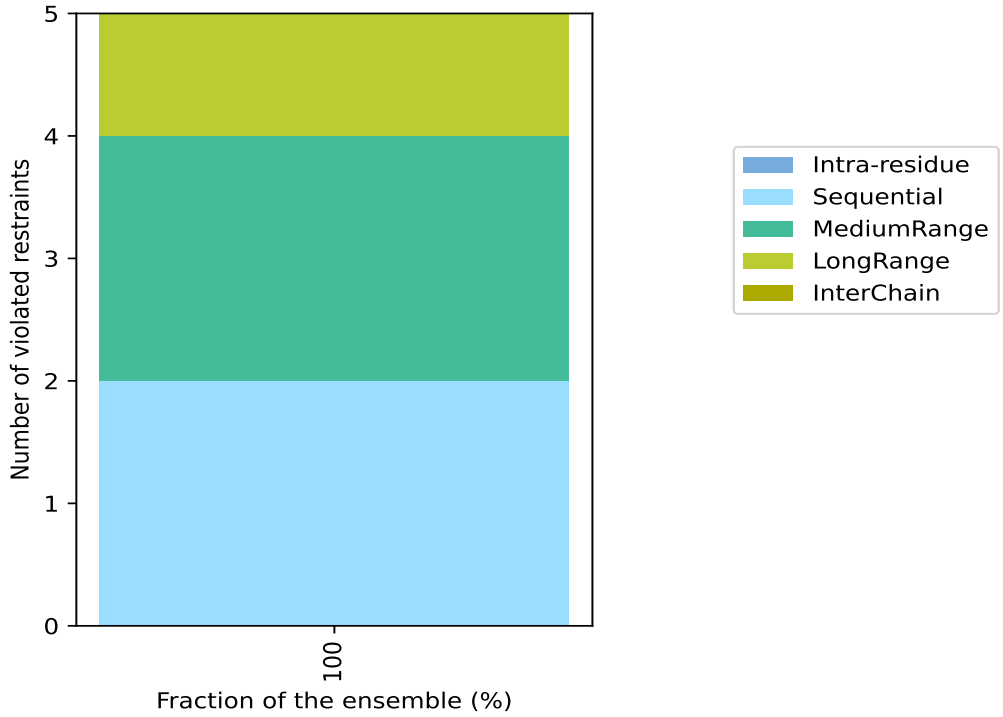
9.3 Distance violation statistics for the ensemble

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, 403(IR:195, SQ:93, MR:53, LR:62, 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]	%
0	2	2	1	0	5	1	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



9.4 Most violated distance restraints in the ensemble

No violations found

9.5 All violated distance restraints

9.5.1 Histogram : Distribution of distance 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

9.5.2 Table : All distance violations

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)
(2,23)	1:A:3:ARG:HE	1:A:5:GLY:HA3	1	5.32

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,103)	1:A:10:VAL:H	1:A:6:PRO:HB3	1	0.4
(2,19)	1:A:3:ARG:HE	1:A:37:GLY:HA2	1	0.39
(2,41)	1:A:4:ILE:H	1:A:3:ARG:HD3	1	0.25
(2,225)	1:A:20:GLU:H	1:A:19:HIS:HB3	1	0.16

10 Dihedral-angle violation analysis

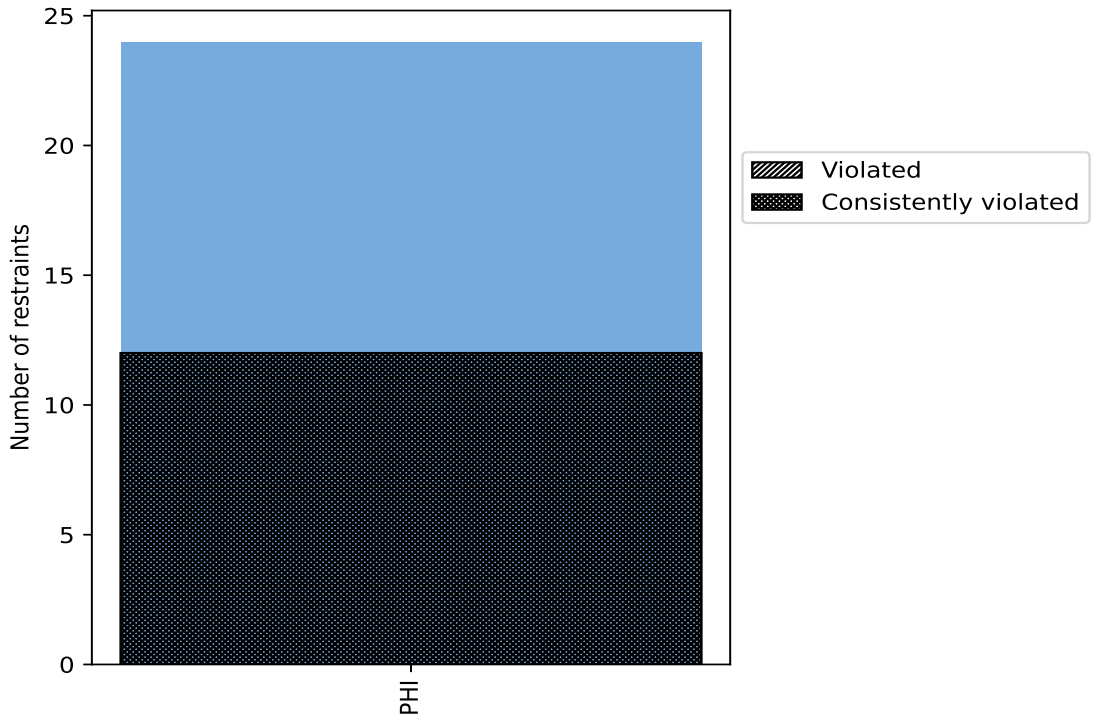
10.1 Summary of dihedral-angle violations

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]
PHI	24	100.0	12	50.0	50.0	12	50.0	50.0
Total	24	100.0	12	50.0	50.0	12	50.0	50.0

[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



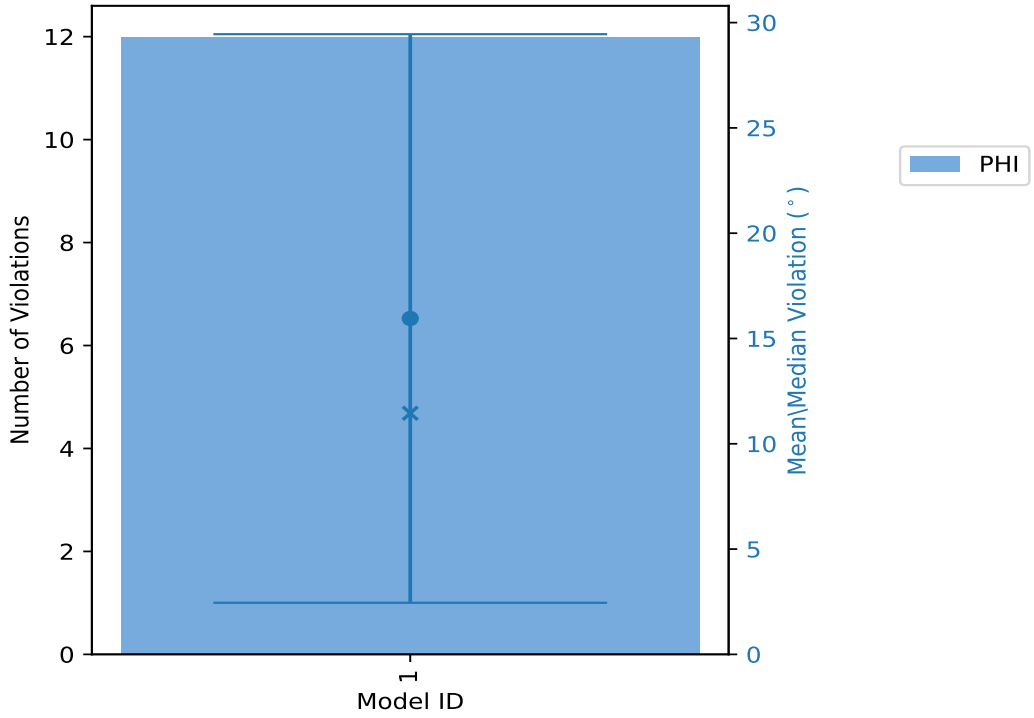
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

10.2 Dihedral-angle violation statistics for each model

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 (°)
	PHI	Total				
1	12	12	15.95	44.6	13.5	11.45

10.2.1 Bar graph : Dihedral violation statistics for each model



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

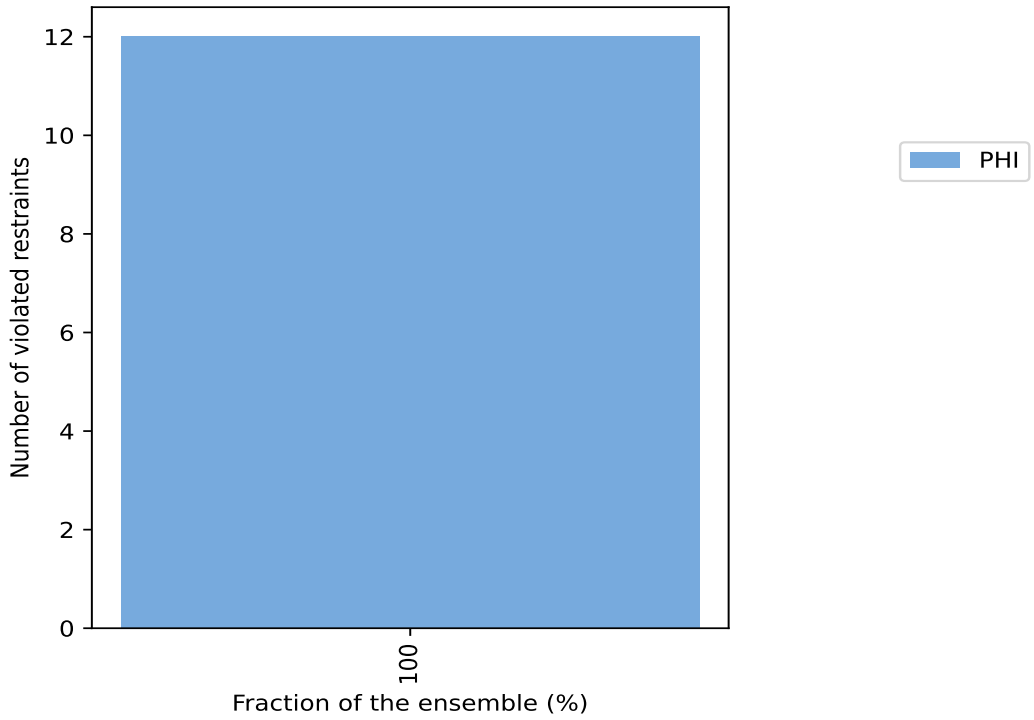
10.3 Dihedral-angle violation statistics for the ensemble

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	
PHI	Total	Count[1]	%
12	12	1	100.0

[1] Number of models with violations

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



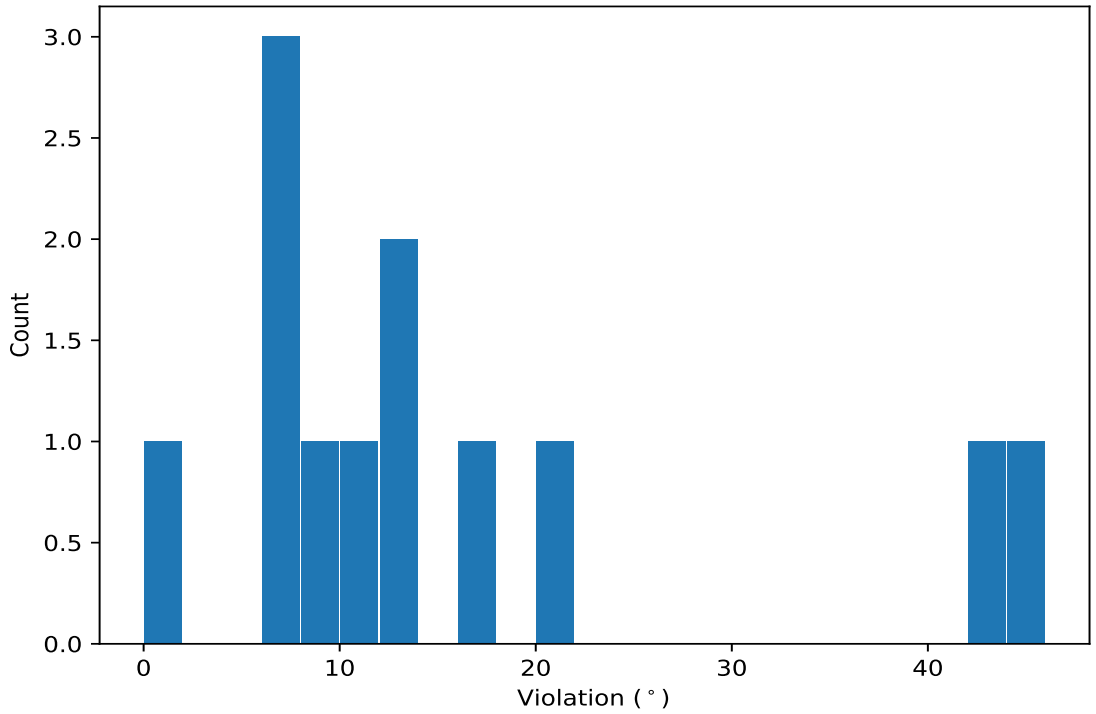
10.4 Most violated dihedral-angle restraints in the ensemble

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.



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 (°)
(1,4)	1:A:6:PRO:C	1:A:7:CYS:N	1:A:7:CYS:CA	1:A:7:CYS:C	1	44.6
(1,2)	1:A:2:VAL:C	1:A:3:ARG:N	1:A:3:ARG:CA	1:A:3:ARG:C	1	43.4
(1,18)	1:A:27:ARG:C	1:A:28:SER:N	1:A:28:SER:CA	1:A:28:SER:C	1	21.3
(1,13)	1:A:21:CYS:C	1:A:22:CYS:N	1:A:22:CYS:CA	1:A:22:CYS:C	1	16.8
(1,10)	1:A:17:GLU:C	1:A:18:ARG:N	1:A:18:ARG:CA	1:A:18:ARG:C	1	13.1
(1,8)	1:A:12:PRO:C	1:A:13:ARG:N	1:A:13:ARG:CA	1:A:13:ARG:C	1	12.0
(1,9)	1:A:14:ILE:C	1:A:15:VAL:N	1:A:15:VAL:CA	1:A:15:VAL:C	1	10.9
(1,6)	1:A:9:GLN:C	1:A:10:VAL:N	1:A:10:VAL:CA	1:A:10:VAL:C	1	8.1
(1,11)	1:A:19:HIS:C	1:A:20:GLU:N	1:A:20:GLU:CA	1:A:20:GLU:C	1	7.1
(1,14)	1:A:22:CYS:C	1:A:23:ARG:N	1:A:23:ARG:CA	1:A:23:ARG:C	1	6.3
(1,22)	1:A:37:GLY:C	1:A:38:MET:N	1:A:38:MET:CA	1:A:38:MET:C	1	6.1
(1,24)	1:A:39:TYR:C	1:A:40:CYS:N	1:A:40:CYS:CA	1:A:40:CYS:C	1	1.7