

Full wwPDB NMR Structure Validation Report (i)

Jun 4, 2023 – 01:18 PM EDT

PDB ID	:	2LI0
BMRB ID	:	17873
Title	:	Mono-O-GalNAc glycosylated Mucin sequence based on MUC2 Mucin glyco-
		protein tandem repeat
Authors	:	Borgert, A.; Heimburg-Molinaro, J.; Lasanajak, Y.; Ju, T.; Liu, M.; Thomp-
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Deposited on	:	2011-08-18

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 (1)) were used in the production of this report:

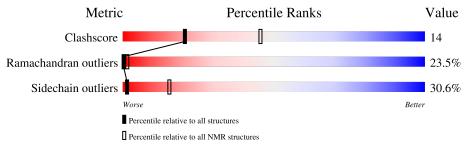
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
BMRB Restraints Analysis	:	v1.2
Ideal geometry (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 91%.

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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f NMR} \ { m archive} \ (\#{ m Entries})$
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	А	9	22%	78%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA and RNA chains that are outliers for geometric criteria:

Mal	Chain	Compound		Total models with violations		
	Ullalli			Chirality	Geometry	
2	А	A2G	10	1	-	



2 Ensemble composition and analysis (i)

This entry contains 62 models.

Cyrange was unable to find well-defined residues.

Error message: The number of core atoms (3) was below the domain threshold value (8).

NmrClust was unable to cluster the ensemble.

Error message: Wrapper check: not enough residues in core to run NmrClust



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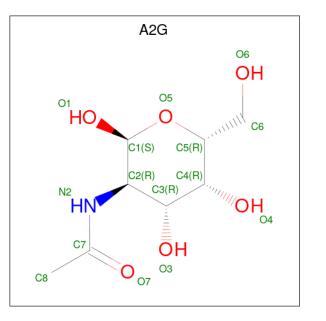
3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 147 atoms, of which 77 are hydrogens and 0 are deuteriums.

• Molecule 1 is a protein called MUC2 Mucin Domain Peptide.

Mol	Chain	Residues	Atoms			Trace		
1	٨	0	Total	С	Η	Ν	0	1
I A	A	A 9	119	36	63	9	11	1

• Molecule 2 is 2-acetamido-2-deoxy-alpha-D-galactopyranose (three-letter code: A2G) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				
2	А	1	Total	С	H	N	0
			28	8	14	1	б



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: MUC2 Mucin Domain Peptide

Chain A:	22%	78%
ACE1 13 14 15 15 15 15 16 15 15 16 16 16 16 16		

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: MUC2 Mucin Domain Peptide



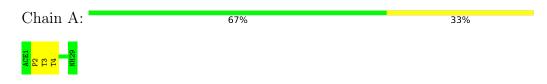
4.2.2 Score per residue for model 2

Chain A:	44%	22%	22%	11%
ACE1 P2 T3 T4 F2 F6 F6 K8 NH29				



4.2.3 Score per residue for model 3

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.4 Score per residue for model 4

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 56% 22% 22%

4.2.5 Score per residue for model 5

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 33% 44% 22%

4.2.6 Score per residue for model 6

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 44% 33% 22%

4.2.7 Score per residue for model 7

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 67% 33%



4.2.8 Score per residue for model 8

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.9 Score per residue for model 9

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A:	44%	33%	22%
ACE1 P2 T5 P6 K8 K8 NH29			

4.2.10 Score per residue for model 10

• Molecule 1: MUC2 Mucin Domain Peptide

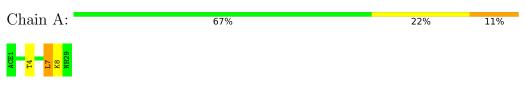
Chain A:	44%	33%	22%
ACE1 P2 T3 T4 F5 F6 K8 NH29 NH29			

4.2.11 Score per residue for model 11

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 44% 11%

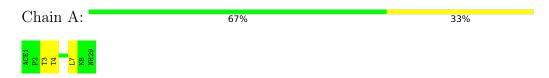
4.2.12 Score per residue for model 12





4.2.13 Score per residue for model 13

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.14 Score per residue for model 14

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 44% 33% 22%

4.2.15 Score per residue for model 15

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 22% 67% 11%

4.2.16 Score per residue for model 16

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 11% 56% 33%

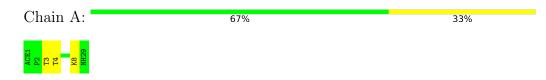
4.2.17 Score per residue for model 17





4.2.18 Score per residue for model 18

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.19 Score per residue for model 19

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 56% 22% 22%

4.2.20 Score per residue for model 20

• Molecule 1: MUC2 Mucin Domain Peptide

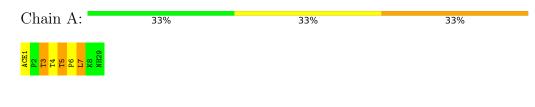
Chain A: 44% 56%

4.2.21 Score per residue for model 21

• Molecule 1: MUC2 Mucin Domain Peptide

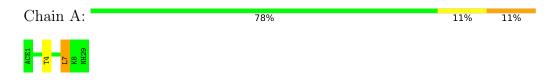
Chain A: 56% 44%

4.2.22 Score per residue for model 22



4.2.23 Score per residue for model 23

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.24 Score per residue for model 24

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 33% 44% 22%

4.2.25 Score per residue for model 25

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A:	22%	67%	11%
ACE 1 22 13 14 14 15 15 17 17 17 17 17 17 17 17 17 17 17 17 17			

4.2.26 Score per residue for model 26

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 56% 22% 22%

4.2.27 Score per residue for model 27





4.2.28 Score per residue for model 28

• Molecule 1: MUC2 Mucin Domain Peptide



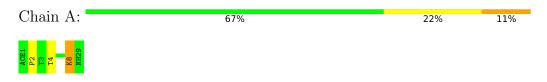
4.2.29 Score per residue for model 29

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A:	78%	11% 11%
ACE1 13 13 113 113 113		

4.2.30 Score per residue for model 30

• Molecule 1: MUC2 Mucin Domain Peptide

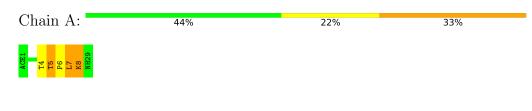


4.2.31 Score per residue for model 31

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 33% 56% 11%

4.2.32 Score per residue for model 32



4.2.33 Score per residue for model 33

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.34 Score per residue for model 34

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 56% 22% 11% 11%

4.2.35 Score per residue for model 35

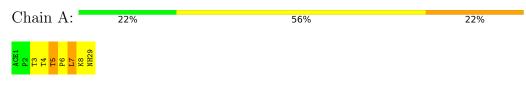
• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 67% 33%

- 4.2.36 Score per residue for model 36
- Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 56% 44%

4.2.37 Score per residue for model 37





4.2.38 Score per residue for model 38

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.39 Score per residue for model 39

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A:	44%	33%	11%	11%
ACE1 P2 T3 T5 T5 T5 T5 M29 M129				

4.2.40 Score per residue for model 40

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A:	56%	33%	11%
ACE1 P2 T3 T4 P6 K8 K8 K8			

11%

4.2.41 Score per residue for model 41

Molecule 1: MUC2 Mucin Domain Peptide
Chain A: 44%

ACE1 P2 T3 T4 L7 K8 NH29

4.2.42 Score per residue for model 42



4.2.43 Score per residue for model 43

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A:	44%	56%
ACE1 P2 T3 T6 P6 MH29		

4.2.44 Score per residue for model 44

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 56% 33% 11%

4.2.45 Score per residue for model 45

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 78% 22%

4.2.46 Score per residue for model 46

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 44% 22% 22% 11%

4.2.47 Score per residue for model 47





4.2.48 Score per residue for model 48

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.49 Score per residue for model 49

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 67% 22% 11%

4.2.50 Score per residue for model 50

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 22% 33% 44%

4.2.51 Score per residue for model 51

Molecule 1: MUC2 Mucin Domain Peptide
Chain A: 67% 33%

4.2.52 Score per residue for model 52





4.2.53 Score per residue for model 53

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A:	56%	22%	22%
ACE1 T5 P6 L7 L7 K8 MH29			

4.2.54 Score per residue for model 54

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 56% 44%

4.2.55 Score per residue for model 55

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 33% 44% 22%

4.2.56 Score per residue for model 56

Molecule 1: MUC2 Mucin Domain Peptide
Chain A: 67% 11% 22%

4.2.57 Score per residue for model 57





4.2.58 Score per residue for model 58

• Molecule 1: MUC2 Mucin Domain Peptide



4.2.59 Score per residue for model 59

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 89% 11%

4.2.60 Score per residue for model 60

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 67% 33%

4.2.61 Score per residue for model 61

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 44% 56%

4.2.62 Score per residue for model 62

• Molecule 1: MUC2 Mucin Domain Peptide

Chain A: 67% 33%



5 Refinement protocol and experimental data overview (i)

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

Of the 100 calculated structures, 62 were deposited, based on the following criterion: structures with the least restraint violations.

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

Software name	Classification	Version
CNS	structure solution	
XPLOR-NIH	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	1
Total number of shifts	99
Number of shifts mapped to atoms	99
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	91%



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: NH2, A2G, ACE $\,$

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	А	56	63	61	2 ± 2
2	А	14	14	12	0±1
All	All	4340	4774	4526	120

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 14.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
Atom-1	Atom-2	Clash(A)	Distance(A)	Worst	Total
1:A:7:LEU:H	1:A:7:LEU:HD13	0.97	1.10	12	1
1:A:5:THR:O	1:A:7:LEU:N	0.92	2.02	53	27
1:A:7:LEU:HD13	1:A:7:LEU:N	0.82	1.88	12	1
1:A:7:LEU:HD13	2:A:10:A2G:O7	0.81	1.73	16	3
1:A:7:LEU:H	1:A:7:LEU:CD1	0.77	1.84	12	3
1:A:7:LEU:HD22	2:A:10:A2G:O7	0.75	1.82	55	1
1:A:7:LEU:HD12	1:A:7:LEU:H	0.74	1.43	23	2
1:A:7:LEU:N	1:A:7:LEU:HD22	0.63	2.09	12	2
1:A:3:THR:O	1:A:3:THR:HG23	0.61	1.94	39	1
1:A:7:LEU:HD22	2:A:10:A2G:H8	0.59	1.72	16	2
1:A:9:NH2:N	2:A:10:A2G:H8	0.59	2.13	37	1

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Continued from pr	Atom-2	$Clash(\lambda)$	Dictorec(Å)	Mo	dels
Atom-1	Atom-2	$\operatorname{Clash}(\operatorname{\AA})$	$\operatorname{Distance}(\operatorname{\AA})$	Worst	Total
1:A:3:THR:HG23	1:A:4:THR:N	0.58	2.13	29	2
1:A:3:THR:HG22	:3:THR:HG22 1:A:4:THR:N		2.13	4	4
1:A:4:THR:HG23	1:A:5:THR:N	0.55	2.17	47	5
1:A:7:LEU:HD23	1:A:7:LEU:N	0.54	2.17	47	2
1:A:7:LEU:N	1:A:7:LEU:CD2	0.52	2.71	6	4
1:A:7:LEU:N	1:A:7:LEU:CD1	0.52	2.71	41	2
1:A:1:ACE:O	1:A:3:THR:N	0.51	2.43	16	5
1:A:6:PRO:O	2:A:10:A2G:C7	0.51	2.59	6	1
1:A:7:LEU:N	1:A:7:LEU:HD12	0.51	2.20	41	1
1:A:7:LEU:HD12	1:A:7:LEU:N	0.50	2.19	23	2
1:A:3:THR:HG22	1:A:4:THR:H	0.50	1.66	4	3
1:A:3:THR:OG1	1:A:4:THR:N	0.50	2.45	1	4
1:A:7:LEU:HD23	2:A:10:A2G:O7	0.49	2.07	10	1
1:A:7:LEU:HD23	1:A:8:LYS:N	0.49	2.21	44	1
1:A:7:LEU:CD1	1:A:7:LEU:O	0.48	2.62	2	1
1:A:7:LEU:HD13	1:A:7:LEU:C	0.47	2.29	52	2
1:A:8:LYS:CG	1:A:9:NH2:N	0.47	2.77	41	1
1:A:7:LEU:CD2	1:A:7:LEU:H	0.47	2.23	40	2
1:A:3:THR:CG2	1:A:4:THR:N	0.47	2.77	4	4
1:A:5:THR:O	1:A:6:PRO:O	0.47	2.33	6	1
1:A:7:LEU:C	1:A:7:LEU:CD1	0.47	2.84	52	2
1:A:7:LEU:CD1	1:A:7:LEU:N	0.46	2.78	57	2
1:A:7:LEU:CD2	2:A:10:A2G:O7	0.46	2.60	55	1
1:A:7:LEU:O	1:A:7:LEU:HD13	0.45	2.11	31	1
1:A:7:LEU:O	1:A:8:LYS:CG	0.45	2.65	36	1
1:A:7:LEU:C	1:A:7:LEU:CD2	0.45	2.85	12	1
1:A:8:LYS:H	1:A:8:LYS:CD	0.45	2.23	30	1
1:A:9:NH2:N	2:A:10:A2G:H8A	0.45	2.27	28	1
1:A:1:ACE:C	1:A:3:THR:N	0.44	2.71	16	5
1:A:7:LEU:CD1	2:A:10:A2G:O7	0.43	2.60	55	1
1:A:3:THR:CG2	1:A:4:THR:H	0.43	2.25	4	2
1:A:8:LYS:O	2:A:10:A2G:O7	0.43	2.37	8	2
1:A:1:ACE:O	1:A:2:PRO:O	0.42	2.38	8	3
1:A:7:LEU:CD2	1:A:8:LYS:H	0.42	2.28	9	1
1:A:9:NH2:N	2:A:10:A2G:O7	0.42	2.53	38	1
1:A:6:PRO:C	1:A:8:LYS:H	0.41	2.19	6	1
1:A:1:ACE:C	1:A:3:THR:H	0.41	2.19	16	1
1:A:7:LEU:HD22	2:A:10:A2G:C8	0.40	2.46	16	1

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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	А	7/9~(78%)	$4\pm1~(54\pm15\%)$	2 ± 1 (23 $\pm13\%$)	$2\pm1 \ (24\pm14\%)$	0 1
All	All	434/558~(78%)	234~(54%)	98~(23%)	102 (24%)	0 1

All 5 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	А	6	PRO	36
1	А	2	PRO	30
1	А	7	LEU	19
1	А	8	LYS	15
1	А	3	THR	2

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	Analysed Rotameric		Perce	entiles
1	А	7/7~(100%)	$5\pm1 (69\pm14\%)$	$2\pm1 (31\pm14\%)$	1	15
All	All	434/434~(100%)	301 (69%)	133 (31%)	1	15

All 5 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	А	4	THR	47
1	А	5	THR	27
1	А	3	THR	23
1	А	8	LYS	19
1	А	7	LEU	17



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)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with |Z| > 2 is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mal	Turne	Chain	Dec	Tiple		ths	
	Type	Chain	nes	LINK	Counts	RMSZ	#Z>2
2	A2G	А	10	1	$14,\!14,\!15$	$0.64{\pm}0.02$	0±0 (0±0%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with |Z| > 2 is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Ма		Chain	Dog	Link		gles	
	Type	Unam	nes	LIIIK	Counts	RMSZ	$\#Z{>}2$
2	A2G	А	10	1	17,19,21	1.05 ± 0.04	$1\pm0~(6\pm1\%)$

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means



no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	A2G	А	10	1	-	$0\pm 0, 6, 23, 26$	$0\pm 0,1,1,1$

There are no bond-length outliers.

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Dec	Type	Atoms	7	$Observed(^{o})$	$Ideal(^{o})$	Moo	dels
IVIOI	Unam	nes	Type	Atoms		Observed()	Ideal(*)	Worst	Total
2	А	10	A2G	C1-O5-C5	3.04	108.08	112.19	47	60
2	А	10	A2G	O5-C1-C2	2.13	107.92	111.29	15	4

All unique chiral outliers are listed below.

Mol	Chain	Res	Type	Atoms	Models (Total)
2	А	10	A2G	C1	1

There are no torsion outliers.

There are no ring outliers.

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 91% for the well-defined parts and 91% 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	99
Number of shifts mapped to atoms	99
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 91%, i.e. 85 atoms were assigned a chemical shift out of a possible 93. 0 out of 1 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^{1}\mathrm{H}$	$^{13}\mathrm{C}$	15 N
Backbone	24/31~(77%)	12/12~(100%)	7/14~(50%)	5/5~(100%)
Sidechain	61/62~(98%)	41/41 (100%)	20/20~(100%)	0/1~(0%)
Overall	85/93~(91%)	53/53~(100%)	27/34~(79%)	5/6 (83%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 91%, i.e. 85 atoms were assigned a chemical shift out of a possible 93. 0 out of 1 assigned methyl groups (LEU and VAL) were assigned stereospecifically.



	Total	$^{1}\mathbf{H}$	$^{13}\mathrm{C}$	$^{15}\mathbf{N}$
Backbone	24/31 (77%)	12/12~(100%)	7/14~(50%)	5/5~(100%)
Sidechain	61/62~(98%)	41/41 (100%)	20/20~(100%)	0/1~(0%)
Overall	85/93~(91%)	53/53~(100%)	27/34~(79%)	5/6~(83%)

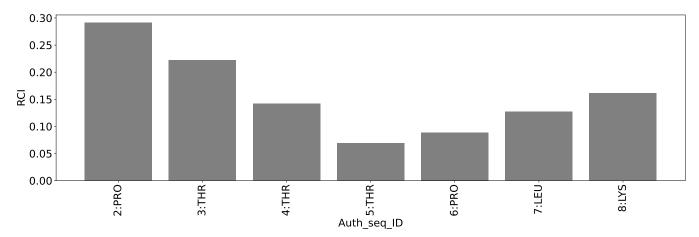
7.1.4 Statistically unusual chemical shifts (i)

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 (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	97
Intra-residue (i-j =0)	65
Sequential (i-j =1)	19
Medium range ($ i-j >1$ and $ i-j <5$)	3
Long range $(i-j \ge 5)$	10
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	7
Number of unmapped restraints	0
Number of restraints per residue	11.6
Number of long range restraints per residue ¹	1.1

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

8.2 Residual restraint violations (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)	0.1	0.15
0.2-0.5 (Medium)	0.1	0.24
>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. There are no dihedral-angle violations



9 Distance violation analysis (i)

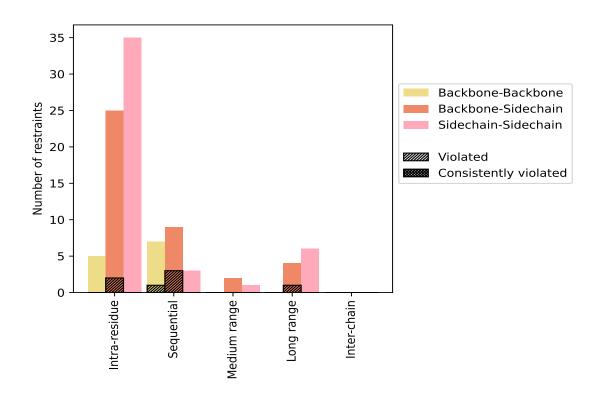
9.1 Summary of distance violations (i)

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

Destructures to me	Count	$\%^1$	Vie	lated ³	3	Consis	tently	Violated ⁴
Restraints type	Count	701	Count	$\%^2$	$\%^1$	Count	$\%^2$	$\%^1$
Intra-residue (i-j =0)	65	67.0	2	3.1	2.1	0	0.0	0.0
Backbone-Backbone	5	5.2	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	25	25.8	2	8.0	2.1	0	0.0	0.0
Sidechain-Sidechain	35	36.1	0	0.0	0.0	0	0.0	0.0
Sequential (i-j =1)	19	19.6	4	21.1	4.1	0	0.0	0.0
Backbone-Backbone	7	7.2	1	14.3	1.0	0	0.0	0.0
Backbone-Sidechain	9	9.3	3	33.3	3.1	0	0.0	0.0
Sidechain-Sidechain	3	3.1	0	0.0	0.0	0	0.0	0.0
Medium range ($ i-j >1 \& i-j <5$)	3	3.1	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	2	2.1	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	1	1.0	0	0.0	0.0	0	0.0	0.0
Long range $(i-j \ge 5)$	10	10.3	1	10.0	1.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	4	4.1	1	25.0	1.0	0	0.0	0.0
Sidechain-Sidechain	6	6.2	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	0	0.0	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	97	100.0	7	7.2	7.2	0	0.0	0.0
Backbone-Backbone	12	12.4	1	8.3	1.0	0	0.0	0.0
Backbone-Sidechain	40	41.2	6	15.0	6.2	0	0.0	0.0
Sidechain-Sidechain	45	46.4	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 (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.

Madal ID		Nun	nber o	f viola	ations	3	Maan (Å)	Mor (Å)	SD^6 (Å)	Madian (Å)
Model ID	IR^{1}	SQ^2	MR^3	LR^4	IC^5	Total	Mean (Å)	Max (Å)	$SD^{*}(A)$	Median (Å)
1	0	0	0	0	0	0	0.0	0.0	0.0	0.0
2	1	0	0	0	0	1	0.24	0.24	0.0	0.24
3	0	0	0	0	0	0	0.0	0.0	0.0	0.0
4	0	0	0	0	0	0	0.0	0.0	0.0	0.0
5	0	0	0	0	0	0	0.0	0.0	0.0	0.0
6	0	0	0	0	0	0	0.0	0.0	0.0	0.0
7	0	0	0	0	0	0	0.0	0.0	0.0	0.0
8	0	0	0	0	0	0	0.0	0.0	0.0	0.0
9	0	0	0	0	0	0	0.0	0.0	0.0	0.0
10	0	0	0	0	0	0	0.0	0.0	0.0	0.0
11	0	0	0	0	0	0	0.0	0.0	0.0	0.0

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Model ID		Nun	nber o	f viola	ations	5	Mass (Å)	Max (Å)	SD^6 (Å)	Madian (8)
Model ID	IR^1	SQ^2	MR^3	LR^4	$ IC^5 $	Total	Mean (Å)	Max (Å)	$SD^{\circ}(A)$	Median (Å)
12	0	0	0	0	0	0	0.0	0.0	0.0	0.0
13	0	0	0	0	0	0	0.0	0.0	0.0	0.0
14	0	0	0	0	0	0	0.0	0.0	0.0	0.0
15	0	0	0	0	0	0	0.0	0.0	0.0	0.0
16	0	0	0	0	0	0	0.0	0.0	0.0	0.0
17	0	0	0	0	0	0	0.0	0.0	0.0	0.0
18	1	0	0	0	0	1	0.13	0.13	0.0	0.13
19	0	0	0	0	0	0	0.0	0.0	0.0	0.0
20	0	0	0	0	0	0	0.0	0.0	0.0	0.0
21	0	0	0	0	0	0	0.0	0.0	0.0	0.0
22	0	0	0	0	0	0	0.0	0.0	0.0	0.0
23	0	0	0	0	0	0	0.0	0.0	0.0	0.0
24	0	0	0	0	0	0	0.0	0.0	0.0	0.0
25	0	0	0	0	0	0	0.0	0.0	0.0	0.0
26	0	0	0	0	0	0	0.0	0.0	0.0	0.0
27	0	0	0	0	0	0	0.0	0.0	0.0	0.0
28	0	0	0	0	0	0	0.0	0.0	0.0	0.0
29	0	0	0	0	0	0	0.0	0.0	0.0	0.0
30	0	0	0	0	0	0	0.0	0.0	0.0	0.0
31	1	0	0	0	0	1	0.23	0.23	0.0	0.23
32	0	0	0	0	0	0	0.0	0.0	0.0	0.0
33	0	0	0	0	0	0	0.0	0.0	0.0	0.0
34	0	0	0	0	0	0	0.0	0.0	0.0	0.0
35	0	0	0	0	0	0	0.0	0.0	0.0	0.0
36	0	0	0	0	0	0	0.0	0.0	0.0	0.0
37	0	0	0	0	0	0	0.0	0.0	0.0	0.0
38	0	0	0	0	0	0	0.0	0.0	0.0	0.0
39	1	0	0	0	0	1	0.24	0.24	0.0	0.24
40	0	0	0	0	0	0	0.0	0.0	0.0	0.0
41	0	0	0	0	0	0	0.0	0.0	0.0	0.0
42	0	0	0	0	0	0	0.0	0.0	0.0	0.0
43	0	0	0	0	0	0	0.0	0.0	0.0	0.0
44	0	2	0	0	0	2	0.14	0.15	0.02	0.14
45	0	0	0	0	0	0	0.0	0.0	0.0	0.0
46	0	0	0	0	0	0	0.0	0.0	0.0	0.0
47	0	0	0	0	0	0	0.0	0.0	0.0	0.0
48	0	0	0	0	0	0	0.0	0.0	0.0	0.0
49	0	1	0	1	0	2	0.15	0.15	0.0	0.15
50	0	0	0	0	0	0	0.0	0.0	0.0	0.0
51	0	0	0	0	0	0	0.0	0.0	0.0	0.0
52	0	0	0	0	0	0	0.0	0.0	0.0	0.0

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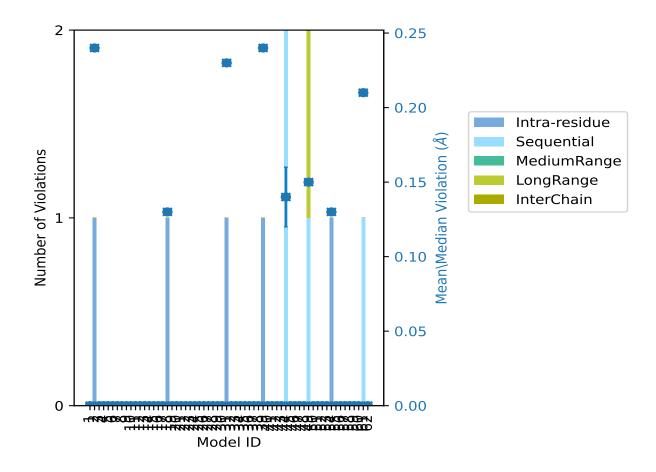


Madal ID		Nun	nber o	f viola	ations	5	Mean (Å)	Max (Å)	SD^6 (Å)	Madian (Å)
Model ID	IR^{1}	SQ^2	MR^3	LR^4	$ IC^5 $	Total	Mean (A)	Max (A)	$SD^{*}(A)$	Median (Å)
53	0	0	0	0	0	0	0.0	0.0	0.0	0.0
54	1	0	0	0	0	1	0.13	0.13	0.0	0.13
55	0	0	0	0	0	0	0.0	0.0	0.0	0.0
56	0	0	0	0	0	0	0.0	0.0	0.0	0.0
57	0	0	0	0	0	0	0.0	0.0	0.0	0.0
58	0	0	0	0	0	0	0.0	0.0	0.0	0.0
59	0	0	0	0	0	0	0.0	0.0	0.0	0.0
60	0	0	0	0	0	0	0.0	0.0	0.0	0.0
61	0	1	0	0	0	1	0.21	0.21	0.0	0.21
62	0	0	0	0	0	0	0.0	0.0	0.0	0.0

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 1 Intra-residue restraints, 2 S
equential 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, 90(IR:63, SQ:15, MR:3, LR:9, IC:0) restraints are not violated in the ensemble.

Nu	mber	of vio	lated	restra	aints	Fractio	n of the ensemble
IR^1	SQ^2	MR^3	LR ⁴	IC ⁵	Total	Count^6	%
0	4	0	1	0	5	1	1.6
1	0	0	0	0	1	2	3.2
1	0	0	0	0	1	3	4.8
0	0	0	0	0	0	4	6.5
0	0	0	0	0	0	5	8.1
0	0	0	0	0	0	6	9.7
0	0	0	0	0	0	7	11.3
0	0	0	0	0	0	8	12.9
0	0	0	0	0	0	9	14.5
0	0	0	0	0	0	10	16.1
0	0	0	0	0	0	11	17.7
0	0	0	0	0	0	12	19.4
0	0	0	0	0	0	13	21.0
0	0	0	0	0	0	14	22.6
0	0	0	0	0	0	15	24.2
0	0	0	0	0	0	16	25.8
0	0	0	0	0	0	17	27.4
0	0	0	0	0	0	18	29.0
0	0	0	0	0	0	19	30.6
0	0	0	0	0	0	20	32.3
0	0	0	0	0	0	21	33.9
0	0	0	0	0	0	22	35.5
0	0	0	0	0	0	23	37.1
0	0	0	0	0	0	24	38.7
0	0	0	0	0	0	25	40.3
0	0	0	0	0	0	26	41.9
0	0	0	0	0	0	27	43.5
0	0	0	0	0	0	28	45.2
0	0	0	0	0	0	29	46.8
0	0	0	0	0	0	30	48.4
0	0	0	0	0	0	31	50.0
0	0	0	0	0	0	32	51.6
0	0	0	0	0	0	33	53.2
0	0	0	0	0	0	34	54.8
0	0	0	0	0	0	35	56.5
0	0	0	0	0	0	36	58.1

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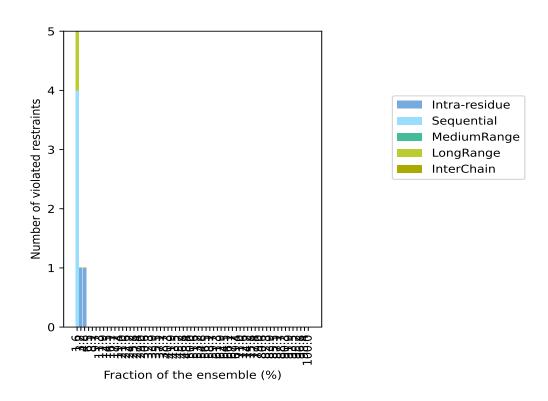


		of vio				Fractio	n of the ensemble
IR^1	SQ^2	MR^3	LR^4	$ IC^5 $	Total	Count^6	%
0	0	0	0	0	0	37	59.7
0	0	0	0	0	0	38	61.3
0	0	0	0	0	0	39	62.9
0	0	0	0	0	0	40	64.5
0	0	0	0	0	0	41	66.1
0	0	0	0	0	0	42	67.7
0	0	0	0	0	0	43	69.4
0	0	0	0	0	0	44	71.0
0	0	0	0	0	0	45	72.6
0	0	0	0	0	0	46	74.2
0	0	0	0	0	0	47	75.8
0	0	0	0	0	0	48	77.4
0	0	0	0	0	0	49	79.0
0	0	0	0	0	0	50	80.6
0	0	0	0	0	0	51	82.3
0	0	0	0	0	0	52	83.9
0	0	0	0	0	0	53	85.5
0	0	0	0	0	0	54	87.1
0	0	0	0	0	0	55	88.7
0	0	0	0	0	0	56	90.3
0	0	0	0	0	0	57	91.9
0	0	0	0	0	0	58	93.5
0	0	0	0	0	0	59	95.2
0	0	0	0	0	0	60	96.8
0	0	0	0	0	0	61	98.4
0	0	0	0	0	0	62	100.0

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 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





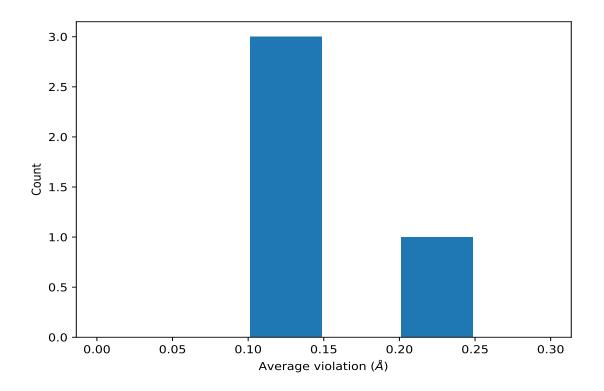
9.3.1 Bar graph : Distance violation statistics for the ensemble (i)

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 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	\mathbf{Models}^1	Mean (Å)	SD^1 (Å)	Median (Å)
(1,85)	1:A:7:LEU:H	1:A:7:LEU:HG	3	0.24	0.0	0.24
(1,25)	1:A:3:THR:H	1:A:3:THR:HG21	2	0.13	0.0	0.13
(1,25)	1:A:3:THR:H	1:A:3:THR:HG22	2	0.13	0.0	0.13
(1,25)	1:A:3:THR:H	1:A:3:THR:HG23	2	0.13	0.0	0.13

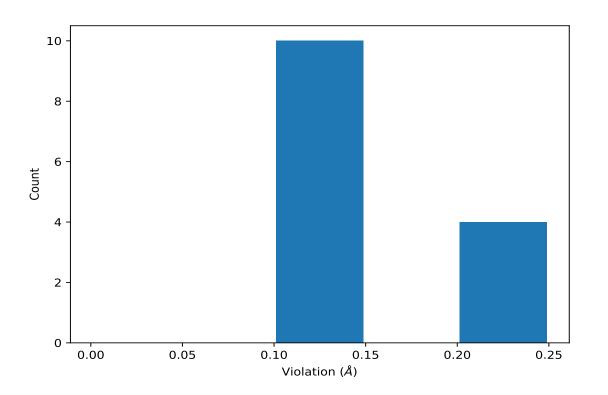
¹Number of violated models, ²Standard deviation

9.5 All violated distance restraints (i)

9.5.1 Histogram : Distribution of distance violations (i)

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





9.5.2 Table : All distance violations (i)

The following table lists the absolute value of the violation for each restraint in the ensemble sorted by its value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,85)	1:A:7:LEU:H	1:A:7:LEU:HG	2	0.24
(1,85)	1:A:7:LEU:H	1:A:7:LEU:HG	39	0.24
(1,85)	1:A:7:LEU:H	1:A:7:LEU:HG	31	0.23
(1,14)	1:A:5:THR:H	1:A:6:PRO:HD3	61	0.21
(1,68)	1:A:4:THR:HA	1:A:5:THR:H	49	0.15
(1,13)	1:A:5:THR:H	1:A:6:PRO:HD2	44	0.15
(1,10)	2:A:10:A2G:HN2	1:A:4:THR:HA	49	0.14
(1,25)	1:A:3:THR:H	1:A:3:THR:HG21	18	0.13
(1,25)	1:A:3:THR:H	1:A:3:THR:HG22	18	0.13
(1,25)	1:A:3:THR:H	1:A:3:THR:HG23	18	0.13
(1,25)	1:A:3:THR:H	1:A:3:THR:HG21	54	0.13
(1,25)	1:A:3:THR:H	1:A:3:THR:HG22	54	0.13
(1,25)	1:A:3:THR:H	1:A:3:THR:HG23	54	0.13
(1,30)	1:A:5:THR:HA	1:A:6:PRO:HD2	44	0.12



10 Dihedral-angle violation analysis (i)

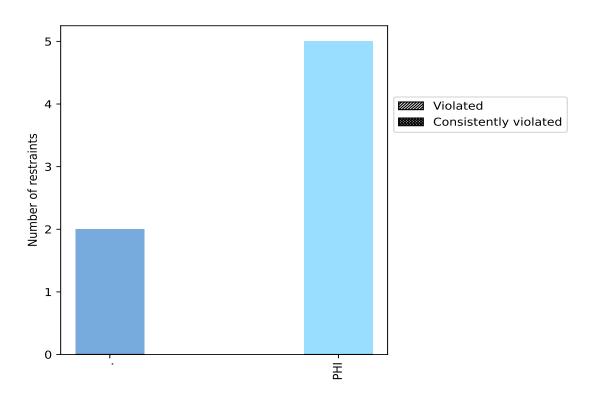
10.1 Summary of dihedral-angle violations (i)

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.

	Count	$\%^1$				Consistently Violated ⁴		
Angle type	Count	70-	Count	$\%^2$	$\%^1$	Count	$\%^2$	$\%^1$
•	2	28.6	0	0.0	0.0	0	0.0	0.0
PHI	5	71.4	0	0.0	0.0	0	0.0	0.0
Total	7	100.0	0	0.0	0.0	0	0.0	0.0

 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	(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)

No violations found

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

No violations found

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

No violations found

10.5 All violated dihedral-angle restraints (i)

No violations found

