



Full wwPDB NMR Structure Validation Report ⓘ

Dec 25, 2024 – 06:17 AM EST

PDB ID : 5YFG
BMRB ID : 36117
Title : SOLUTION STRUCTURE OF HUMAN MOG1
Authors : Hu, Q.; Liu, Y.; Bao, X.; Liu, H.
Deposited on : 2017-09-21

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 ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
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) : 2.40

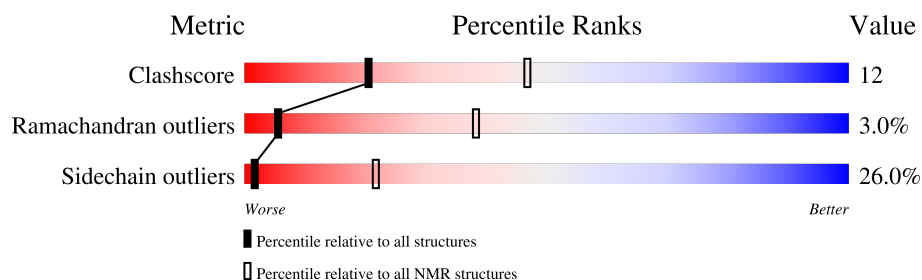
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 66%.

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	210492	14027
Ramachandran outliers	207382	12486
Sidechain outliers	206894	12463

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	194	

2 Ensemble composition and analysis

This entry contains 20 models. Model 17 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:5-A:112, A:119-A:145, A:162-A:180 (154)	0.79	17

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

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

The models can be grouped into 4 clusters. No single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2858 atoms, of which 1418 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Ran guanine nucleotide release factor.

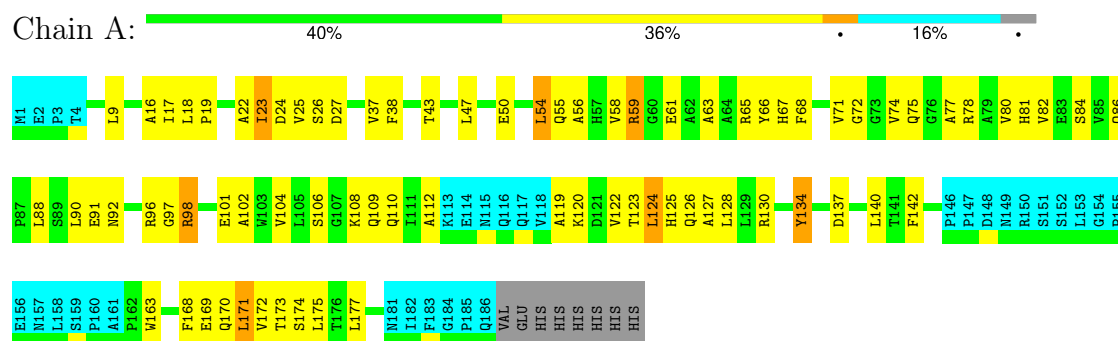
Mol	Chain	Residues	Atoms						Trace
1	A	186	Total	C	H	N	O	S	0
			2858	907	1418	253	275	5	

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	187	VAL	-	expression tag	UNP Q9HD47
A	188	GLU	-	expression tag	UNP Q9HD47
A	189	HIS	-	expression tag	UNP Q9HD47
A	190	HIS	-	expression tag	UNP Q9HD47
A	191	HIS	-	expression tag	UNP Q9HD47
A	192	HIS	-	expression tag	UNP Q9HD47
A	193	HIS	-	expression tag	UNP Q9HD47
A	194	HIS	-	expression tag	UNP Q9HD47

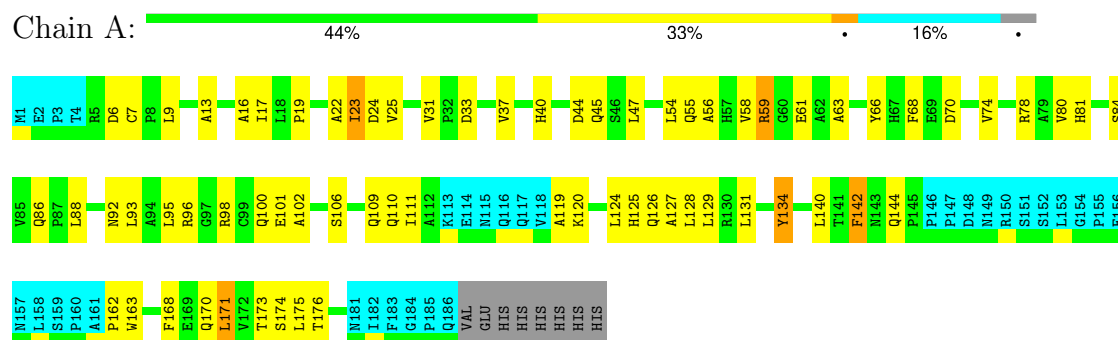
4.2.2 Score per residue for model 2

- Molecule 1: Ran guanine nucleotide release factor



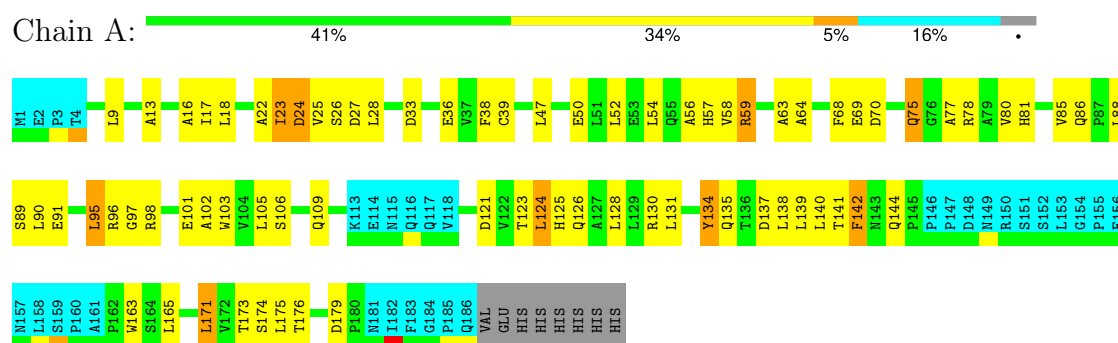
4.2.3 Score per residue for model 3

- Molecule 1: Ran guanine nucleotide release factor



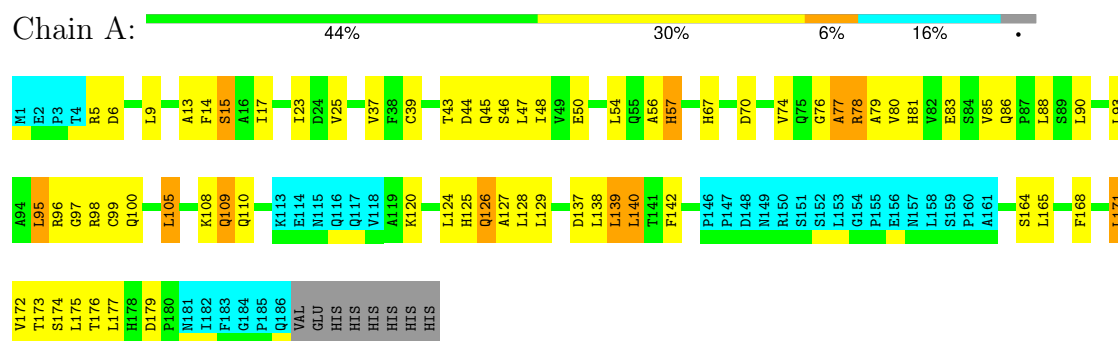
4.2.4 Score per residue for model 4

- Molecule 1: Ran guanine nucleotide release factor



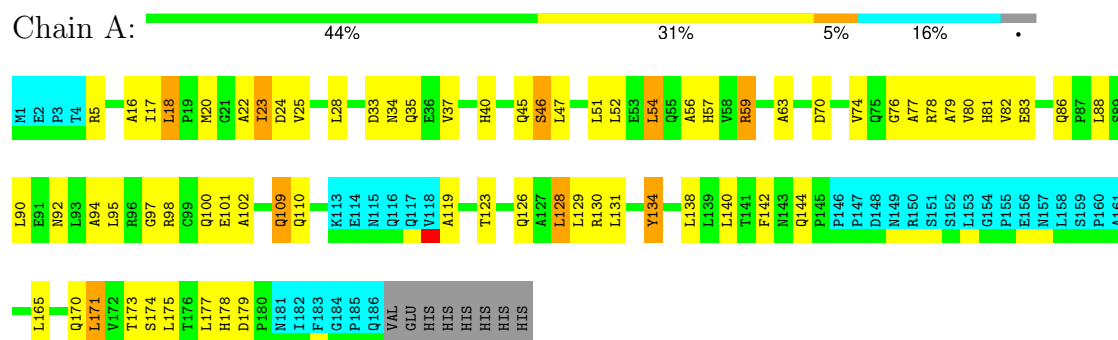
4.2.5 Score per residue for model 5

- Molecule 1: Ran guanine nucleotide release factor



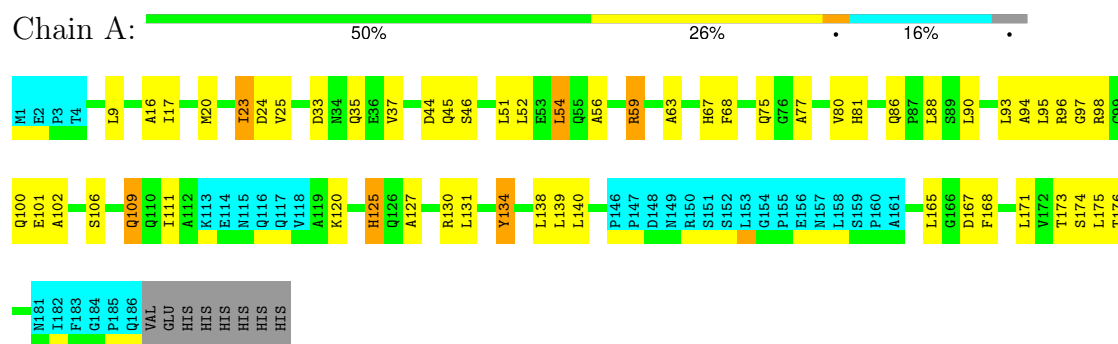
4.2.6 Score per residue for model 6

- Molecule 1: Ran guanine nucleotide release factor



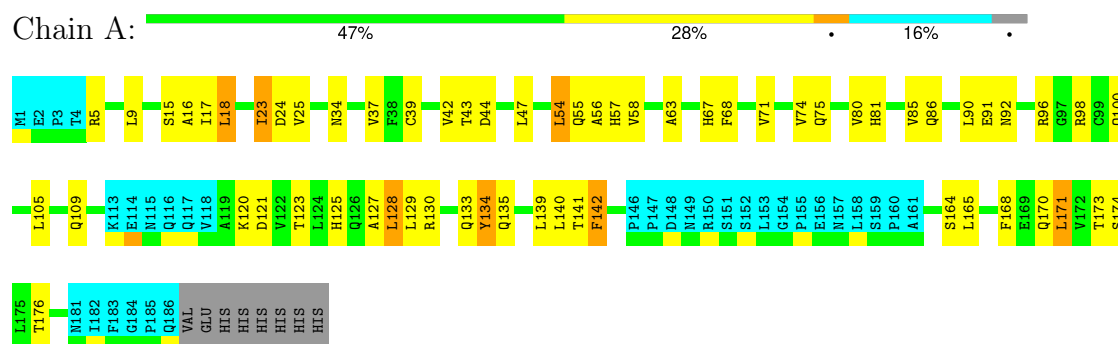
4.2.7 Score per residue for model 7

- Molecule 1: Ran guanine nucleotide release factor



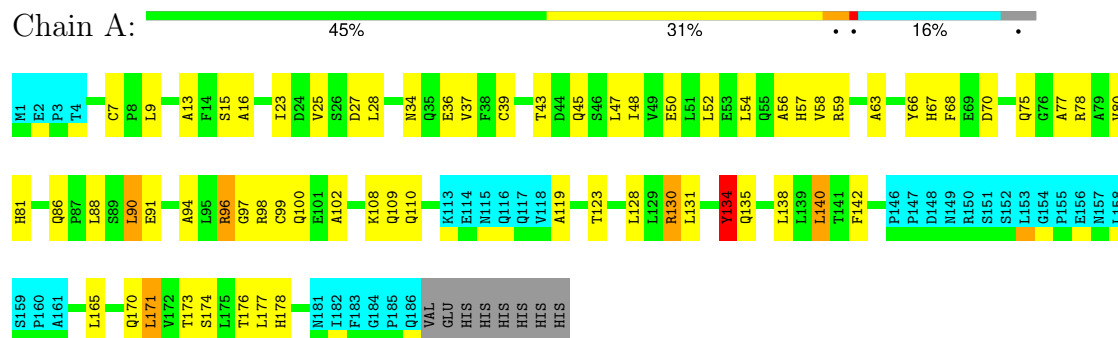
4.2.8 Score per residue for model 8

- Molecule 1: Ran guanine nucleotide release factor



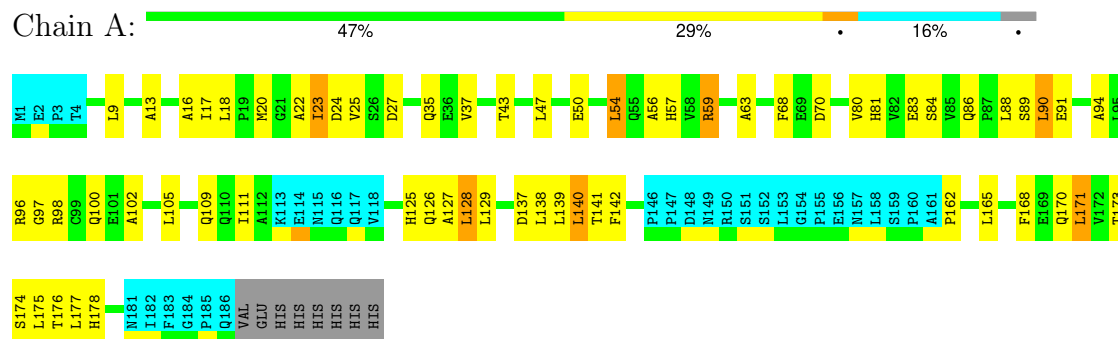
4.2.9 Score per residue for model 9

- Molecule 1: Ran guanine nucleotide release factor



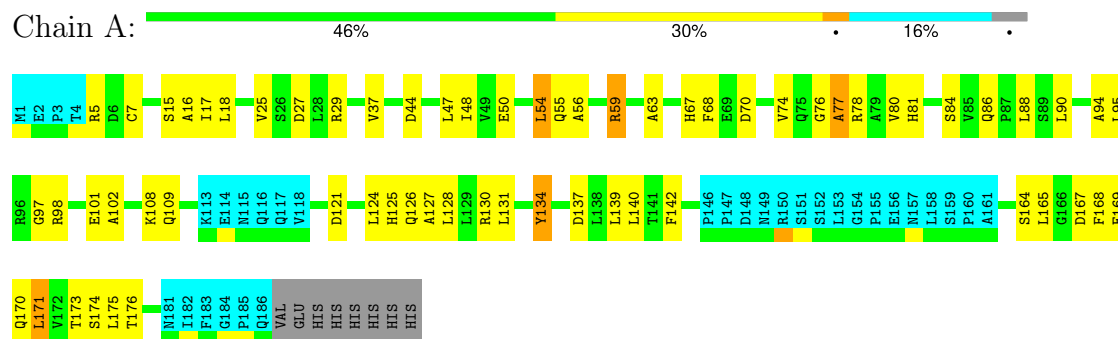
4.2.10 Score per residue for model 10

- Molecule 1: Ran guanine nucleotide release factor



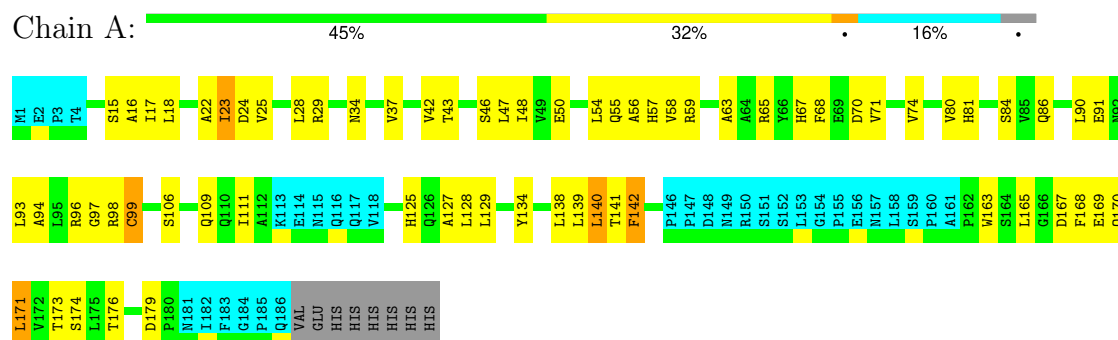
4.2.11 Score per residue for model 11

- Molecule 1: Ran guanine nucleotide release factor



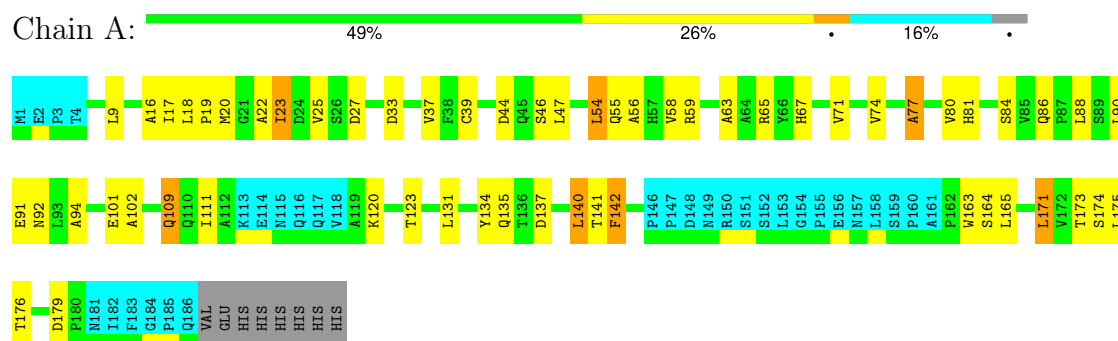
4.2.12 Score per residue for model 12

- Molecule 1: Ran guanine nucleotide release factor



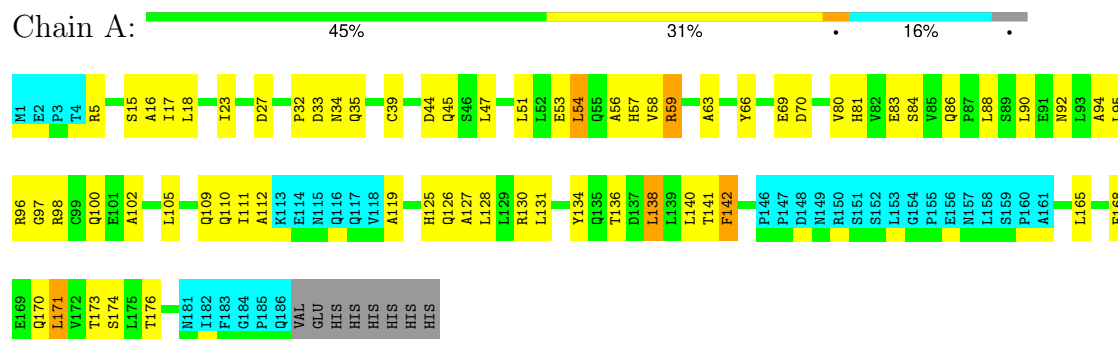
4.2.13 Score per residue for model 13

- Molecule 1: Ran guanine nucleotide release factor



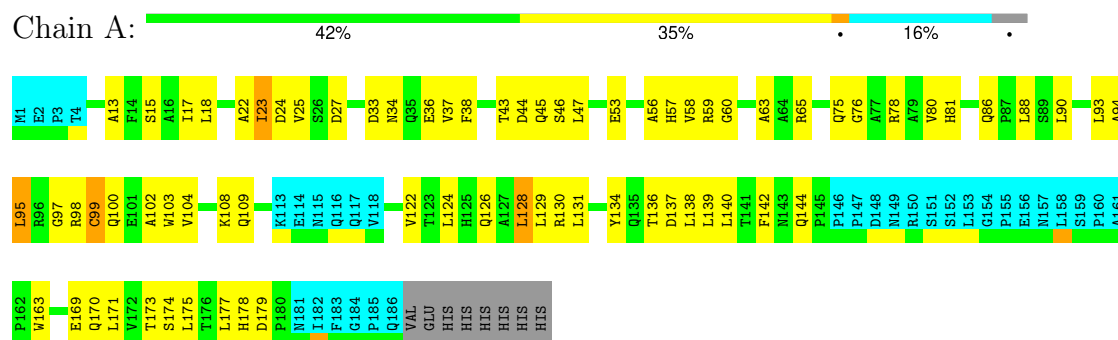
4.2.14 Score per residue for model 14

- Molecule 1: Ran guanine nucleotide release factor



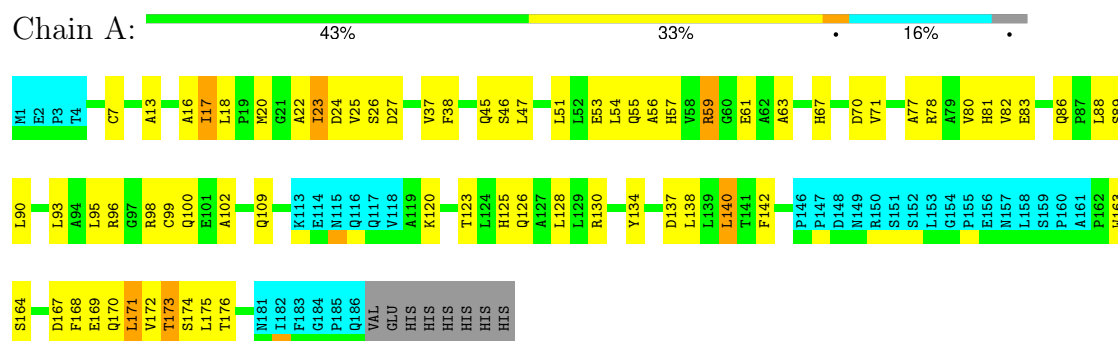
4.2.15 Score per residue for model 15

- Molecule 1: Ran guanine nucleotide release factor



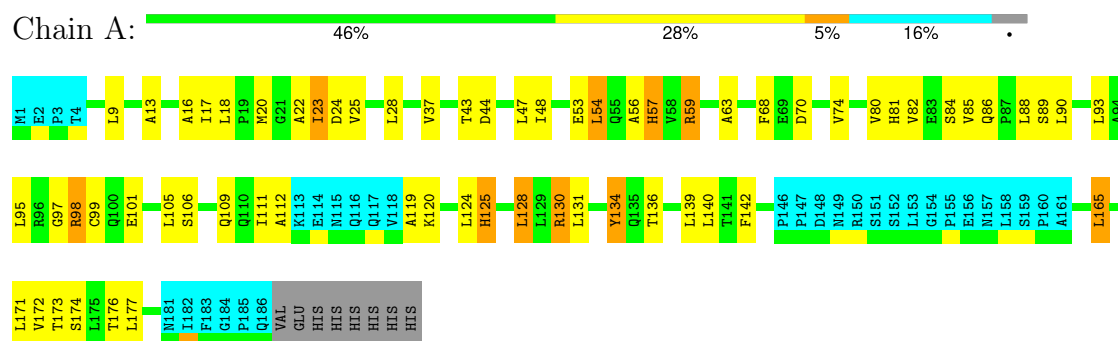
4.2.16 Score per residue for model 16

- Molecule 1: Ran guanine nucleotide release factor



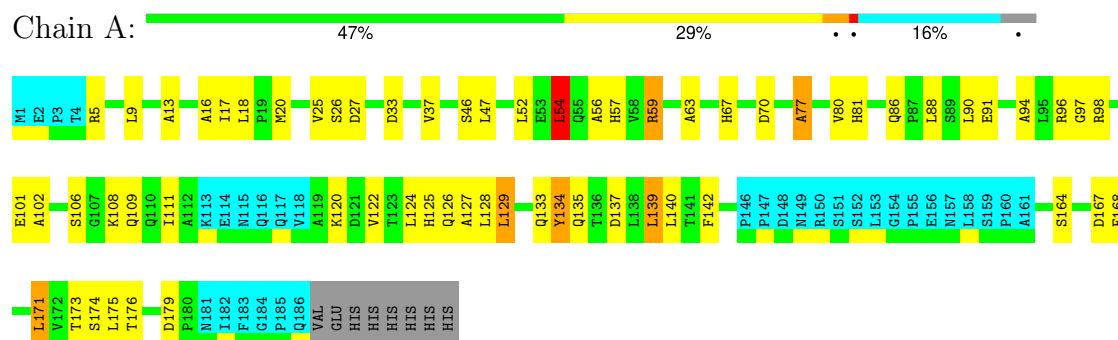
4.2.17 Score per residue for model 17 (medoid)

- Molecule 1: Ran guanine nucleotide release factor



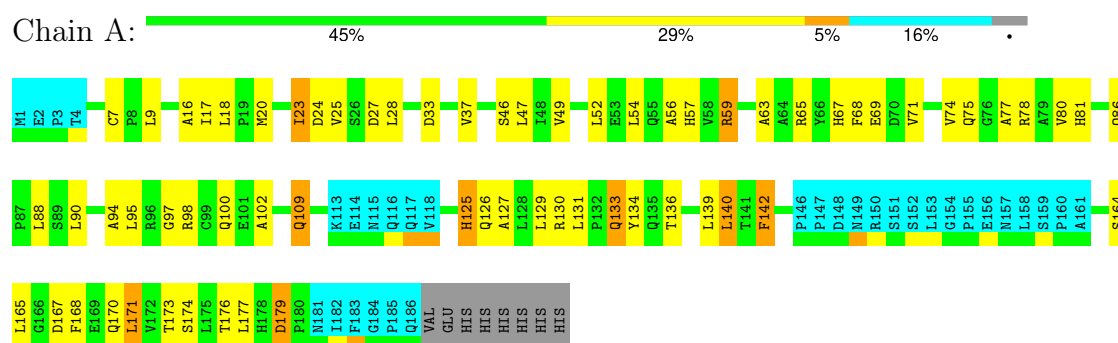
4.2.18 Score per residue for model 18

- Molecule 1: Ran guanine nucleotide release factor



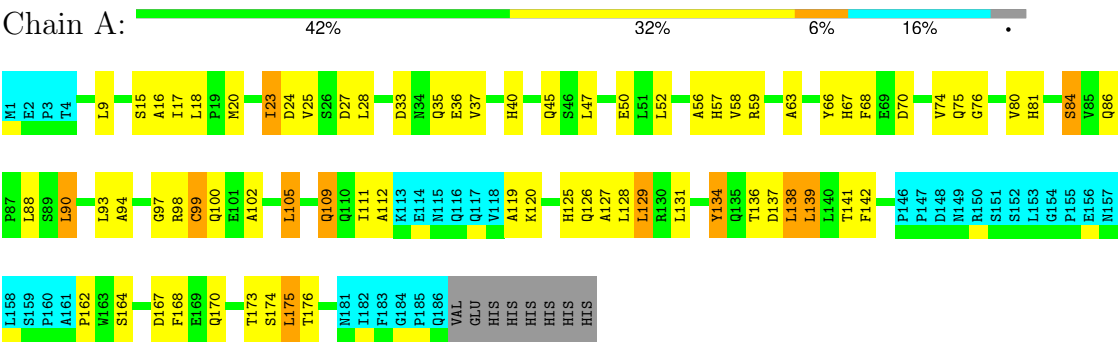
4.2.19 Score per residue for model 19

- Molecule 1: Ran guanine nucleotide release factor



4.2.20 Score per residue for model 20

- Molecule 1: Ran guanine nucleotide release factor



5 Refinement protocol and experimental data overview

The models were refined using the following method: *SIMULATED ANNEALING, DISTANCE GEOMETRY, 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	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	1638
Number of shifts mapped to atoms	1638
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	66%

6 Model quality

6.1 Standard geometry

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

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	1196	1185	1179	28±3
All	All	23920	23700	23580	563

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 12.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:138:LEU:HD22	1:A:175:LEU:HD21	0.97	1.34	1	1
1:A:112:ALA:HB2	1:A:119:ALA:HB2	0.94	1.38	17	4
1:A:140:LEU:HD21	1:A:171:LEU:HD12	0.93	1.37	10	5
1:A:129:LEU:HD11	1:A:175:LEU:HD13	0.72	1.58	3	1
1:A:88:LEU:HB2	1:A:102:ALA:HB3	0.72	1.61	2	8
1:A:140:LEU:CD1	1:A:171:LEU:HD12	0.71	2.16	2	1
1:A:19:PRO:HD2	1:A:22:ALA:HB2	0.70	1.61	3	3
1:A:177:LEU:C	1:A:177:LEU:HD13	0.70	2.07	9	4
1:A:127:ALA:HB2	1:A:168:PHE:CZ	0.70	2.21	3	11
1:A:47:LEU:HD21	1:A:171:LEU:HG	0.70	1.63	4	2
1:A:95:LEU:HD22	1:A:176:THR:HB	0.69	1.64	16	2
1:A:47:LEU:HD13	1:A:142:PHE:CD2	0.69	2.21	4	2
1:A:58:VAL:HG11	1:A:66:TYR:CD2	0.69	2.23	9	3
1:A:80:VAL:HG22	1:A:109:GLN:CB	0.67	2.19	9	18

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:127:ALA:HB3	1:A:140:LEU:HD22	0.67	1.66	2	1
1:A:47:LEU:HD11	1:A:142:PHE:CZ	0.66	2.26	14	6
1:A:59:ARG:HA	1:A:63:ALA:HB2	0.65	1.69	20	15
1:A:37:VAL:HG12	1:A:48:ILE:HG12	0.65	1.67	12	1
1:A:68:PHE:CE1	1:A:80:VAL:HG11	0.64	2.27	9	12
1:A:140:LEU:HD23	1:A:171:LEU:HD12	0.64	1.68	6	1
1:A:88:LEU:HD21	1:A:172:VAL:HG11	0.64	1.69	17	1
1:A:58:VAL:HG12	1:A:63:ALA:HA	0.64	1.69	1	5
1:A:74:VAL:HG13	1:A:75:GLN:HG3	0.62	1.72	20	1
1:A:140:LEU:CD2	1:A:171:LEU:HD12	0.62	2.24	8	5
1:A:88:LEU:HB3	1:A:102:ALA:HB3	0.62	1.71	19	8
1:A:17:ILE:HD12	1:A:174:SER:HA	0.62	1.69	5	3
1:A:52:LEU:HD13	1:A:139:LEU:HD13	0.61	1.72	7	1
1:A:140:LEU:HD21	1:A:171:LEU:HB3	0.61	1.70	12	1
1:A:128:LEU:HD21	1:A:139:LEU:HD12	0.61	1.71	15	1
1:A:99:CYS:SG	1:A:129:LEU:HD11	0.61	2.35	20	1
1:A:37:VAL:HG13	1:A:48:ILE:CD1	0.61	2.25	11	2
1:A:173:THR:O	1:A:176:THR:HG23	0.60	1.95	12	15
1:A:18:LEU:HD12	1:A:22:ALA:CB	0.60	2.27	17	5
1:A:124:LEU:C	1:A:124:LEU:HD22	0.60	2.17	4	1
1:A:23:ILE:HD13	1:A:24:ASP:N	0.60	2.11	8	13
1:A:140:LEU:HD13	1:A:171:LEU:HB3	0.59	1.73	3	2
1:A:67:HIS:O	1:A:71:VAL:HG23	0.59	1.97	8	7
1:A:88:LEU:CD2	1:A:172:VAL:HG11	0.59	2.28	1	2
1:A:56:ALA:O	1:A:58:VAL:HG23	0.59	1.98	8	1
1:A:47:LEU:HD21	1:A:142:PHE:CE2	0.58	2.33	13	1
1:A:110:GLN:HG3	1:A:119:ALA:HB1	0.58	1.74	9	3
1:A:77:ALA:HB1	1:A:109:GLN:HB2	0.58	1.76	19	6
1:A:95:LEU:HB2	1:A:176:THR:HG23	0.58	1.73	4	1
1:A:70:ASP:O	1:A:74:VAL:HG23	0.58	1.98	5	2
1:A:79:ALA:HB3	1:A:110:GLN:HB2	0.58	1.75	6	1
1:A:54:LEU:HD21	1:A:59:ARG:HD3	0.58	1.74	1	2
1:A:138:LEU:CD2	1:A:175:LEU:HD21	0.57	2.21	1	1
1:A:94:ALA:H	1:A:173:THR:HG22	0.57	1.60	12	12
1:A:88:LEU:HD22	1:A:92:ASN:ND2	0.57	2.15	6	1
1:A:25:VAL:HG22	1:A:37:VAL:O	0.56	2.00	5	18
1:A:68:PHE:CE1	1:A:80:VAL:HG21	0.56	2.34	1	1
1:A:54:LEU:HD11	1:A:59:ARG:HG2	0.56	1.76	6	1
1:A:63:ALA:HB3	1:A:128:LEU:CD2	0.56	2.31	10	2
1:A:68:PHE:HE1	1:A:80:VAL:HG11	0.56	1.60	4	9
1:A:177:LEU:HD13	1:A:178:HIS:N	0.56	2.14	15	4

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:168:PHE:O	1:A:172:VAL:HG23	0.56	2.01	5	2
1:A:127:ALA:HB2	1:A:168:PHE:HZ	0.56	1.60	3	8
1:A:95:LEU:HD23	1:A:178:HIS:CE1	0.56	2.35	15	1
1:A:98:ARG:HD3	1:A:177:LEU:HD23	0.55	1.76	17	1
1:A:138:LEU:HD11	1:A:175:LEU:HD22	0.55	1.77	20	1
1:A:54:LEU:HD11	1:A:130:ARG:HG3	0.55	1.75	17	1
1:A:54:LEU:HD11	1:A:59:ARG:HD3	0.55	1.77	13	1
1:A:110:GLN:HG3	1:A:119:ALA:HB3	0.55	1.77	14	1
1:A:80:VAL:HG22	1:A:109:GLN:HB3	0.55	1.79	17	5
1:A:16:ALA:HA	1:A:174:SER:O	0.55	2.02	16	18
1:A:127:ALA:HB3	1:A:140:LEU:HB2	0.55	1.76	12	3
1:A:125:HIS:CE1	1:A:165:LEU:HD21	0.55	2.36	17	1
1:A:59:ARG:CA	1:A:63:ALA:HB2	0.54	2.31	7	13
1:A:47:LEU:HD13	1:A:142:PHE:CZ	0.54	2.37	18	1
1:A:37:VAL:HG22	1:A:48:ILE:HG23	0.54	1.78	17	2
1:A:80:VAL:HG22	1:A:109:GLN:HB2	0.54	1.80	7	6
1:A:90:LEU:HD22	1:A:96:ARG:HD2	0.54	1.79	9	1
1:A:140:LEU:HD11	1:A:171:LEU:HD12	0.53	1.79	2	2
1:A:54:LEU:HD21	1:A:59:ARG:HB3	0.53	1.80	12	3
1:A:138:LEU:HG	1:A:175:LEU:HD22	0.53	1.79	16	1
1:A:37:VAL:HG13	1:A:48:ILE:HG12	0.53	1.79	5	2
1:A:129:LEU:HD12	1:A:138:LEU:CB	0.53	2.33	15	1
1:A:63:ALA:HB3	1:A:128:LEU:HG	0.53	1.80	6	3
1:A:7:CYS:SG	1:A:18:LEU:HD23	0.53	2.44	19	2
1:A:49:VAL:CG2	1:A:171:LEU:HD11	0.53	2.33	19	1
1:A:138:LEU:HD22	1:A:175:LEU:CD1	0.52	2.34	5	1
1:A:67:HIS:CD2	1:A:139:LEU:HD11	0.52	2.39	11	1
1:A:47:LEU:HD21	1:A:171:LEU:CG	0.52	2.33	4	1
1:A:122:VAL:HG13	1:A:122:VAL:O	0.52	2.05	18	1
1:A:17:ILE:C	1:A:18:LEU:HD22	0.52	2.25	14	9
1:A:124:LEU:HD21	1:A:126:GLN:NE2	0.52	2.20	5	1
1:A:52:LEU:HD11	1:A:139:LEU:HD12	0.52	1.82	20	1
1:A:47:LEU:HD11	1:A:142:PHE:CE2	0.51	2.40	5	8
1:A:177:LEU:C	1:A:177:LEU:CD1	0.51	2.79	9	4
1:A:76:GLY:O	1:A:77:ALA:HB2	0.51	2.05	11	1
1:A:68:PHE:CD2	1:A:105:LEU:HD22	0.51	2.40	20	1
1:A:18:LEU:N	1:A:18:LEU:HD22	0.51	2.20	18	1
1:A:23:ILE:HG22	1:A:39:CYS:O	0.51	2.04	5	5
1:A:88:LEU:CB	1:A:102:ALA:HB3	0.51	2.35	4	3
1:A:80:VAL:HG13	1:A:108:LYS:O	0.51	2.06	2	2
1:A:94:ALA:HB2	1:A:173:THR:HB	0.50	1.82	15	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:111:ILE:C	1:A:111:ILE:HD13	0.50	2.25	1	1
1:A:127:ALA:HB3	1:A:140:LEU:HD21	0.50	1.83	19	1
1:A:140:LEU:HD13	1:A:171:LEU:HD12	0.50	1.83	19	2
1:A:68:PHE:CZ	1:A:80:VAL:HG21	0.50	2.42	1	1
1:A:72:GLY:O	1:A:77:ALA:HB3	0.50	2.07	1	1
1:A:177:LEU:HD22	1:A:178:HIS:N	0.50	2.22	10	4
1:A:25:VAL:CG2	1:A:37:VAL:HG23	0.50	2.37	15	2
1:A:17:ILE:HG21	1:A:170:GLN:NE2	0.49	2.22	19	2
1:A:95:LEU:HD22	1:A:176:THR:HG22	0.49	1.84	14	1
1:A:19:PRO:CD	1:A:22:ALA:HB2	0.49	2.38	13	1
1:A:140:LEU:HD22	1:A:171:LEU:HD12	0.48	1.83	3	2
1:A:67:HIS:CG	1:A:139:LEU:HD21	0.48	2.43	5	2
1:A:17:ILE:C	1:A:18:LEU:HD23	0.48	2.29	6	3
1:A:177:LEU:HD22	1:A:178:HIS:H	0.48	1.67	15	3
1:A:74:VAL:O	1:A:74:VAL:HG12	0.48	2.08	13	8
1:A:177:LEU:HD11	1:A:179:ASP:O	0.48	2.08	19	1
1:A:85:VAL:HG12	1:A:105:LEU:HA	0.48	1.85	17	3
1:A:122:VAL:HG13	1:A:124:LEU:HD11	0.48	1.86	2	2
1:A:77:ALA:HB1	1:A:109:GLN:CG	0.48	2.39	5	1
1:A:140:LEU:HD21	1:A:171:LEU:CD1	0.48	2.26	10	3
1:A:80:VAL:HG23	1:A:108:LYS:O	0.47	2.09	1	1
1:A:51:LEU:O	1:A:52:LEU:HD22	0.47	2.09	6	1
1:A:58:VAL:HG21	1:A:66:TYR:CD2	0.47	2.44	20	2
1:A:17:ILE:HD13	1:A:170:GLN:NE2	0.47	2.24	3	1
1:A:63:ALA:HB3	1:A:128:LEU:HD21	0.47	1.84	10	1
1:A:18:LEU:HD12	1:A:22:ALA:HB3	0.47	1.85	16	2
1:A:64:ALA:HB2	1:A:103:TRP:CD1	0.47	2.45	4	1
1:A:47:LEU:HD22	1:A:142:PHE:CE2	0.47	2.44	12	1
1:A:32:PRO:HG2	1:A:35:GLN:HB2	0.47	1.86	14	1
1:A:129:LEU:HD12	1:A:138:LEU:HB2	0.47	1.86	15	1
1:A:131:LEU:HD12	1:A:131:LEU:N	0.47	2.25	6	7
1:A:54:LEU:HB2	1:A:135:GLN:HB3	0.47	1.86	8	2
1:A:40:HIS:CE1	1:A:47:LEU:HD13	0.47	2.45	20	1
1:A:37:VAL:HG13	1:A:48:ILE:CG1	0.47	2.39	5	1
1:A:47:LEU:HD23	1:A:171:LEU:HG	0.47	1.87	11	2
1:A:90:LEU:HD23	1:A:93:LEU:HD13	0.47	1.86	5	1
1:A:17:ILE:HD13	1:A:17:ILE:N	0.47	2.25	17	3
1:A:129:LEU:HD22	1:A:129:LEU:N	0.47	2.24	5	2
1:A:47:LEU:HD21	1:A:142:PHE:CZ	0.47	2.45	8	2
1:A:25:VAL:HG21	1:A:37:VAL:CG1	0.47	2.40	18	1
1:A:93:LEU:HD21	1:A:99:CYS:SG	0.46	2.50	12	2

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:7:CYS:O	1:A:9:LEU:HD12	0.46	2.10	3	1
1:A:94:ALA:HB2	1:A:173:THR:CG2	0.46	2.41	9	1
1:A:88:LEU:HD13	1:A:172:VAL:HG11	0.46	1.87	2	1
1:A:47:LEU:HD21	1:A:171:LEU:CD2	0.46	2.41	12	1
1:A:47:LEU:HD11	1:A:142:PHE:CE1	0.46	2.46	13	1
1:A:111:ILE:O	1:A:111:ILE:HG23	0.46	2.11	1	9
1:A:90:LEU:HD23	1:A:96:ARG:HG2	0.46	1.86	10	2
1:A:131:LEU:HB2	1:A:136:THR:CG2	0.45	2.41	20	6
1:A:95:LEU:HD13	1:A:177:LEU:O	0.45	2.11	6	1
1:A:138:LEU:HD22	1:A:175:LEU:HD13	0.45	1.87	5	1
1:A:128:LEU:CD2	1:A:139:LEU:HD12	0.45	2.41	8	2
1:A:54:LEU:HD11	1:A:130:ARG:CG	0.45	2.40	17	1
1:A:47:LEU:HD11	1:A:142:PHE:CD2	0.45	2.46	3	2
1:A:7:CYS:SG	1:A:16:ALA:HB3	0.45	2.51	16	1
1:A:22:ALA:HB1	1:A:38:PHE:CD2	0.45	2.47	2	2
1:A:47:LEU:HD23	1:A:171:LEU:CD2	0.45	2.41	5	1
1:A:78:ARG:O	1:A:80:VAL:HG23	0.45	2.12	5	1
1:A:127:ALA:HB3	1:A:140:LEU:CD2	0.44	2.41	19	1
1:A:22:ALA:HB1	1:A:38:PHE:CD1	0.44	2.47	4	1
1:A:93:LEU:HD11	1:A:99:CYS:SG	0.44	2.52	16	2
1:A:65:ARG:NH1	1:A:105:LEU:HD12	0.44	2.27	1	1
1:A:31:VAL:O	1:A:31:VAL:HG13	0.44	2.13	3	1
1:A:173:THR:O	1:A:174:SER:C	0.44	2.56	5	1
1:A:52:LEU:HD13	1:A:67:HIS:CE1	0.44	2.47	20	1
1:A:54:LEU:HD13	1:A:135:GLN:HG3	0.44	1.88	13	1
1:A:129:LEU:HD21	1:A:140:LEU:HD13	0.44	1.89	18	1
1:A:67:HIS:CD2	1:A:128:LEU:HD11	0.44	2.47	20	1
1:A:54:LEU:HD21	1:A:59:ARG:CD	0.44	2.41	1	1
1:A:58:VAL:HG12	1:A:63:ALA:CA	0.44	2.42	4	1
1:A:67:HIS:CG	1:A:139:LEU:HD11	0.44	2.47	18	1
1:A:47:LEU:HD12	1:A:141:THR:O	0.44	2.12	4	1
1:A:15:SER:O	1:A:175:LEU:HA	0.44	2.13	5	2
1:A:19:PRO:HG2	1:A:22:ALA:HB2	0.44	1.89	13	1
1:A:60:GLY:HA2	1:A:103:TRP:CD1	0.44	2.48	15	1
1:A:40:HIS:CB	1:A:163:TRP:CH2	0.43	3.01	3	1
1:A:51:LEU:HD13	1:A:138:LEU:HG	0.43	1.89	7	1
1:A:171:LEU:HD13	1:A:171:LEU:O	0.43	2.13	1	3
1:A:82:VAL:O	1:A:82:VAL:HG23	0.43	2.14	17	3
1:A:139:LEU:C	1:A:140:LEU:HD12	0.43	2.34	18	1
1:A:95:LEU:HD12	1:A:176:THR:HB	0.43	1.90	5	1
1:A:59:ARG:N	1:A:63:ALA:HB2	0.43	2.28	7	2

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:171:LEU:O	1:A:174:SER:HB2	0.43	2.14	5	4
1:A:25:VAL:HG22	1:A:25:VAL:O	0.43	2.14	4	1
1:A:15:SER:O	1:A:174:SER:O	0.43	2.37	15	2
1:A:140:LEU:CB	1:A:171:LEU:HD12	0.43	2.44	3	1
1:A:90:LEU:HD11	1:A:99:CYS:O	0.43	2.13	20	1
1:A:52:LEU:CD1	1:A:139:LEU:HD12	0.42	2.44	18	1
1:A:52:LEU:HD23	1:A:67:HIS:CE1	0.42	2.48	9	1
1:A:52:LEU:HD22	1:A:67:HIS:CE1	0.42	2.49	19	1
1:A:127:ALA:CB	1:A:140:LEU:HD22	0.42	2.42	2	1
1:A:85:VAL:HG12	1:A:105:LEU:HG	0.42	1.91	5	1
1:A:82:VAL:HG23	1:A:82:VAL:O	0.42	2.15	16	1
1:A:88:LEU:HD12	1:A:102:ALA:CB	0.42	2.45	16	1
1:A:88:LEU:HD12	1:A:102:ALA:HB1	0.42	1.90	16	1
1:A:128:LEU:HD23	1:A:139:LEU:HD12	0.42	1.91	17	1
1:A:40:HIS:HB3	1:A:163:TRP:CH2	0.42	2.50	3	1
1:A:63:ALA:HB3	1:A:130:ARG:NH1	0.42	2.30	9	1
1:A:111:ILE:HG23	1:A:120:LYS:HD2	0.42	1.92	13	1
1:A:104:VAL:HG13	1:A:168:PHE:CZ	0.41	2.50	2	1
1:A:25:VAL:HG23	1:A:37:VAL:HG23	0.41	1.92	6	2
1:A:58:VAL:HG11	1:A:66:TYR:CB	0.41	2.45	1	1
1:A:35:GLN:HG3	1:A:50:GLU:HG3	0.41	1.91	10	1
1:A:47:LEU:CD2	1:A:171:LEU:HD23	0.41	2.45	3	1
1:A:47:LEU:HD13	1:A:142:PHE:CE1	0.41	2.50	18	1
1:A:18:LEU:HD22	1:A:18:LEU:N	0.41	2.30	20	1
1:A:58:VAL:HG21	1:A:66:TYR:CE2	0.41	2.49	20	1
1:A:54:LEU:HD11	1:A:59:ARG:CD	0.41	2.46	13	1
1:A:14:PHE:CE1	1:A:177:LEU:HD13	0.41	2.50	5	1
1:A:51:LEU:C	1:A:52:LEU:HD12	0.41	2.36	7	1
1:A:51:LEU:HD12	1:A:138:LEU:CD2	0.41	2.44	14	1
1:A:129:LEU:HD11	1:A:175:LEU:CD1	0.41	2.40	3	1
1:A:55:GLN:O	1:A:58:VAL:HG23	0.41	2.16	13	1
1:A:18:LEU:HD11	1:A:38:PHE:CD2	0.41	2.50	16	1
1:A:125:HIS:HB3	1:A:168:PHE:CE1	0.41	2.50	19	1
1:A:98:ARG:HG2	1:A:98:ARG:O	0.41	2.16	2	1
1:A:140:LEU:HD13	1:A:171:LEU:CB	0.41	2.43	3	1
1:A:46:SER:O	1:A:47:LEU:HD12	0.41	2.15	6	1
1:A:40:HIS:HE1	1:A:47:LEU:HD13	0.41	1.75	6	1
1:A:125:HIS:HB3	1:A:168:PHE:CE2	0.41	2.51	7	1
1:A:72:GLY:HA2	1:A:77:ALA:HB3	0.40	1.93	2	1
1:A:79:ALA:N	1:A:110:GLN:HB3	0.40	2.31	5	1
1:A:88:LEU:HD13	1:A:104:VAL:HG23	0.40	1.93	15	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:56:ALA:O	1:A:57:HIS:CG	0.40	2.74	8	1
1:A:23:ILE:HD13	1:A:23:ILE:C	0.40	2.37	17	1
1:A:54:LEU:HD21	1:A:59:ARG:NH1	0.40	2.31	7	1
1:A:134:TYR:CD1	1:A:134:TYR:N	0.40	2.89	9	1
1:A:54:LEU:HD11	1:A:59:ARG:HD2	0.40	1.94	10	1
1:A:55:GLN:O	1:A:58:VAL:HG22	0.40	2.16	12	1
1:A:94:ALA:N	1:A:173:THR:HG22	0.40	2.31	15	1
1:A:124:LEU:HD12	1:A:124:LEU:C	0.40	2.36	18	1
1:A:49:VAL:HG22	1:A:171:LEU:HD11	0.40	1.93	19	1

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	154/194 (79%)	134±2 (87±1%)	15±2 (10±1%)	5±1 (3±1%)	5	38
All	All	3080/3880 (79%)	2684 (87%)	305 (10%)	91 (3%)	5	38

All 12 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	56	ALA	19
1	A	134	TYR	17
1	A	97	GLY	16
1	A	57	HIS	14
1	A	77	ALA	9
1	A	75	GLN	5
1	A	76	GLY	4
1	A	54	LEU	3
1	A	19	PRO	1
1	A	23	ILE	1
1	A	84	SER	1
1	A	162	PRO	1

6.3.2 Protein sidechains ⓘ

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	130/167 (78%)	96±4 (74±3%)	34±4 (26±3%)	2	22
All	All	2600/3340 (78%)	1925 (74%)	675 (26%)	2	22

All 89 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	81	HIS	20
1	A	86	GLN	20
1	A	98	ARG	19
1	A	171	LEU	19
1	A	90	LEU	17
1	A	54	LEU	16
1	A	125	HIS	16
1	A	59	ARG	15
1	A	126	GLN	14
1	A	165	LEU	14
1	A	128	LEU	14
1	A	9	LEU	13
1	A	100	GLN	13
1	A	130	ARG	13
1	A	23	ILE	13
1	A	27	ASP	12
1	A	142	PHE	12
1	A	137	ASP	11
1	A	134	TYR	11
1	A	70	ASP	11
1	A	44	ASP	10
1	A	84	SER	10
1	A	101	GLU	10
1	A	78	ARG	10
1	A	96	ARG	10
1	A	170	GLN	10
1	A	33	ASP	10
1	A	43	THR	9
1	A	120	LYS	9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Models (Total)
1	A	175	LEU	9
1	A	45	GLN	9
1	A	46	SER	9
1	A	140	LEU	9
1	A	20	MET	9
1	A	95	LEU	8
1	A	138	LEU	8
1	A	139	LEU	8
1	A	91	GLU	8
1	A	179	ASP	8
1	A	164	SER	8
1	A	141	THR	7
1	A	163	TRP	7
1	A	50	GLU	7
1	A	106	SER	7
1	A	123	THR	7
1	A	28	LEU	7
1	A	15	SER	7
1	A	167	ASP	7
1	A	5	ARG	6
1	A	99	CYS	6
1	A	109	GLN	6
1	A	34	ASN	6
1	A	129	LEU	6
1	A	26	SER	5
1	A	53	GLU	5
1	A	108	LYS	5
1	A	144	GLN	5
1	A	55	GLN	5
1	A	65	ARG	5
1	A	92	ASN	5
1	A	124	LEU	5
1	A	169	GLU	5
1	A	83	GLU	5
1	A	35	GLN	4
1	A	133	GLN	4
1	A	36	GLU	4
1	A	89	SER	4
1	A	105	LEU	4
1	A	18	LEU	3
1	A	29	ARG	3
1	A	61	GLU	3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Models (Total)
1	A	173	THR	3
1	A	93	LEU	3
1	A	69	GLU	3
1	A	75	GLN	3
1	A	121	ASP	3
1	A	51	LEU	2
1	A	47	LEU	2
1	A	6	ASP	2
1	A	135	GLN	2
1	A	57	HIS	2
1	A	42	VAL	2
1	A	7	CYS	2
1	A	17	ILE	2
1	A	111	ILE	1
1	A	24	ASP	1
1	A	39	CYS	1
1	A	52	LEU	1
1	A	88	LEU	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 oligosaccharides 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 ⓘ

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 66% for the well-defined parts and 65% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *mog1-shifts_gaiHB1_HB3.txt*

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	1638
Number of shifts mapped to atoms	1638
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	6

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	172	-0.18 ± 0.08	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	151	-0.13 ± 0.16	None needed (< 0.5 ppm)
$^{13}\text{C}'$	152	0.28 ± 0.09	None needed (< 0.5 ppm)
^{15}N	160	-0.28 ± 0.37	None needed (< 0.5 ppm)

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 66%, i.e. 1391 atoms were assigned a chemical shift out of a possible 2098. 0 out of 37 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	697/760 (92%)	286/308 (93%)	273/308 (89%)	138/144 (96%)
Sidechain	694/1188 (58%)	494/778 (63%)	200/368 (54%)	0/42 (0%)

Continued on next page...

Continued from previous page...

	Total	¹ H	¹³ C	¹⁵ N
Aromatic	0/150 (0%)	0/74 (0%)	0/62 (0%)	0/14 (0%)
Overall	1391/2098 (66%)	780/1160 (67%)	473/738 (64%)	138/200 (69%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 65%, i.e. 1637 atoms were assigned a chemical shift out of a possible 2507. 0 out of 40 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	817/910 (90%)	333/368 (90%)	324/372 (87%)	160/170 (94%)
Sidechain	820/1437 (57%)	575/937 (61%)	245/447 (55%)	0/53 (0%)
Aromatic	0/160 (0%)	0/79 (0%)	0/67 (0%)	0/14 (0%)
Overall	1637/2507 (65%)	908/1384 (66%)	569/886 (64%)	160/237 (68%)

7.1.4 Statistically unusual chemical shifts [i](#)

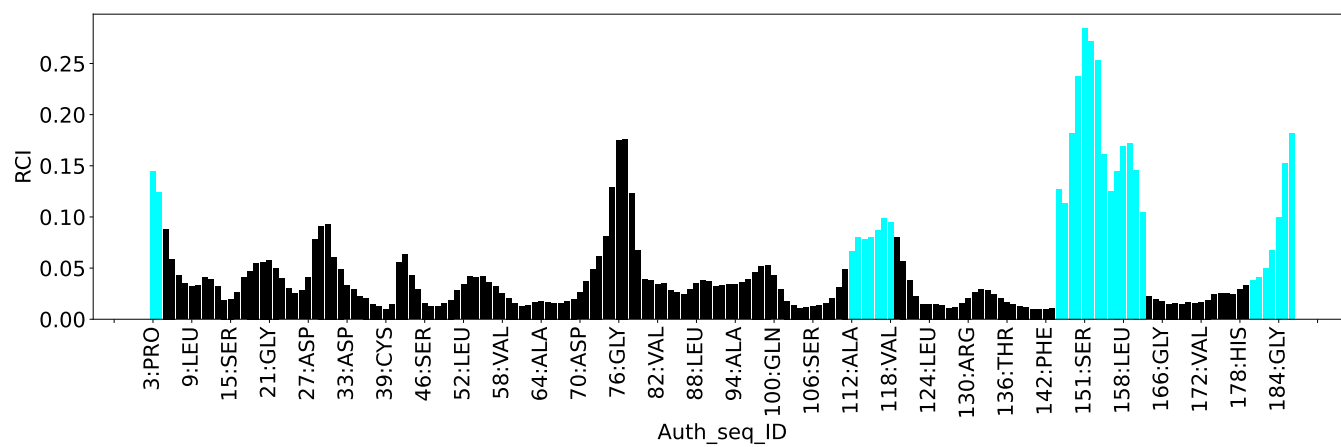
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	123	THR	CB	29.62	61.12 – 78.27	-23.4
1	A	96	ARG	CD	23.59	38.57 – 47.75	-21.3
1	A	17	ILE	CG2	37.26	10.93 – 24.12	15.0
1	A	74	VAL	CG1	41.21	14.71 – 28.29	14.5
1	A	74	VAL	CG2	41.21	13.71 – 28.88	13.1
1	A	17	ILE	CD1	32.92	5.18 – 21.60	11.9

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	2632
Intra-residue ($ i-j =0$)	759
Sequential ($ i-j =1$)	756
Medium range ($ i-j >1$ and $ i-j <5$)	308
Long range ($ i-j \geq 5$)	719
Inter-chain	0
Hydrogen bond restraints	90
Disulfide bond restraints	0
Total dihedral-angle restraints	194
Number of unmapped restraints	0
Number of restraints per residue	14.6
Number of long range restraints per residue ¹	4.1

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

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)	3.7	0.16
0.2-0.5 (Medium)	0.2	0.31
>0.5 (Large)	1.1	2.3

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)	1.1	3.97
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

9 Distance violation analysis ⓘ

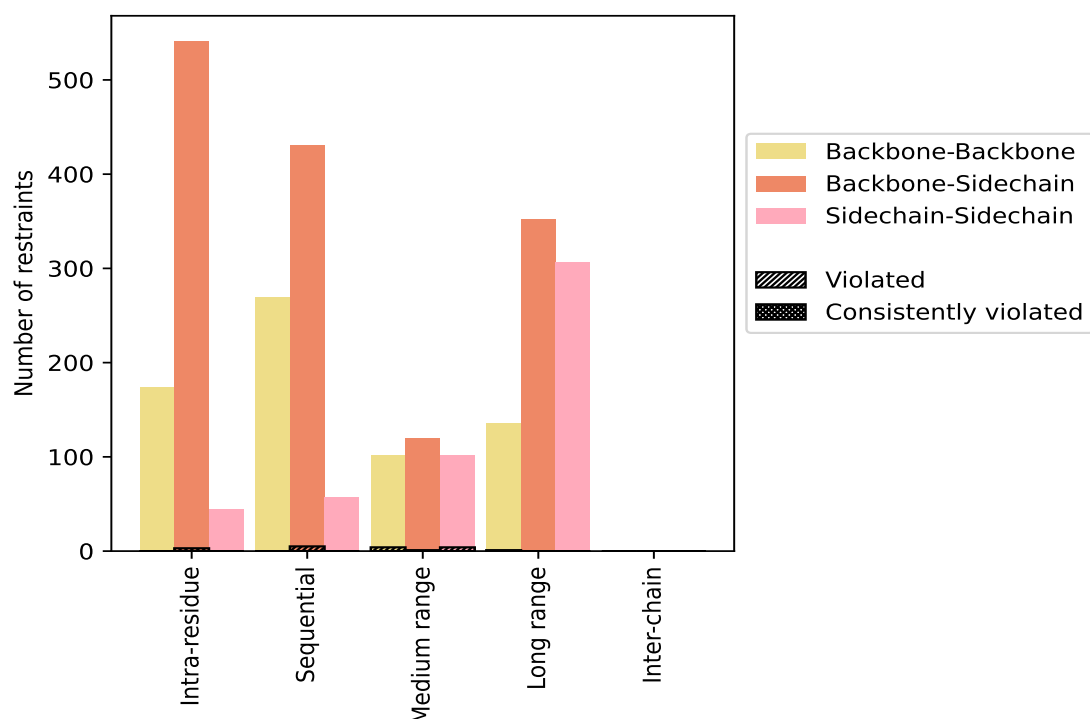
9.1 Summary of distance violations ⓘ

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	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	759	28.8	3	0.4	0.1	0	0.0	0.0
Backbone-Backbone	174	6.6	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	541	20.6	3	0.6	0.1	0	0.0	0.0
Sidechain-Sidechain	44	1.7	0	0.0	0.0	0	0.0	0.0
Sequential ($i-j =1$)	756	28.7	5	0.7	0.2	0	0.0	0.0
Backbone-Backbone	269	10.2	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	430	16.3	5	1.2	0.2	0	0.0	0.0
Sidechain-Sidechain	57	2.2	0	0.0	0.0	0	0.0	0.0
Medium range ($i-j >1$ & $i-j <5$)	308	11.7	9	2.9	0.3	0	0.0	0.0
Backbone-Backbone	102	3.9	4	3.9	0.2	0	0.0	0.0
Backbone-Sidechain	104	4.0	1	1.0	0.0	0	0.0	0.0
Sidechain-Sidechain	102	3.9	4	3.9	0.2	0	0.0	0.0
Long range ($i-j \geq 5$)	719	27.3	1	0.1	0.0	0	0.0	0.0
Backbone-Backbone	135	5.1	1	0.7	0.0	0	0.0	0.0
Backbone-Sidechain	278	10.6	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	306	11.6	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	90	3.4	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	2632	100.0	18	0.7	0.7	0	0.0	0.0
Backbone-Backbone	680	25.8	5	0.7	0.2	0	0.0	0.0
Backbone-Sidechain	1443	54.8	9	0.6	0.3	0	0.0	0.0
Sidechain-Sidechain	509	19.3	4	0.8	0.2	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.

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	1	0	4	1	0	6	0.63	1.99	0.74	0.13
2	0	0	2	0	0	2	0.15	0.15	0.0	0.15
3	1	0	5	1	0	7	0.52	1.31	0.48	0.23
4	1	1	2	1	0	5	0.13	0.15	0.02	0.12
5	0	1	5	0	0	6	0.21	0.67	0.21	0.12
6	1	3	4	1	0	9	0.41	1.71	0.53	0.14
7	1	2	2	1	0	6	0.13	0.15	0.02	0.13
8	0	1	5	0	0	6	0.68	1.58	0.56	0.56
9	0	2	2	1	0	5	0.13	0.15	0.01	0.13
10	1	2	2	0	0	5	0.13	0.15	0.02	0.13

Continued on next page...

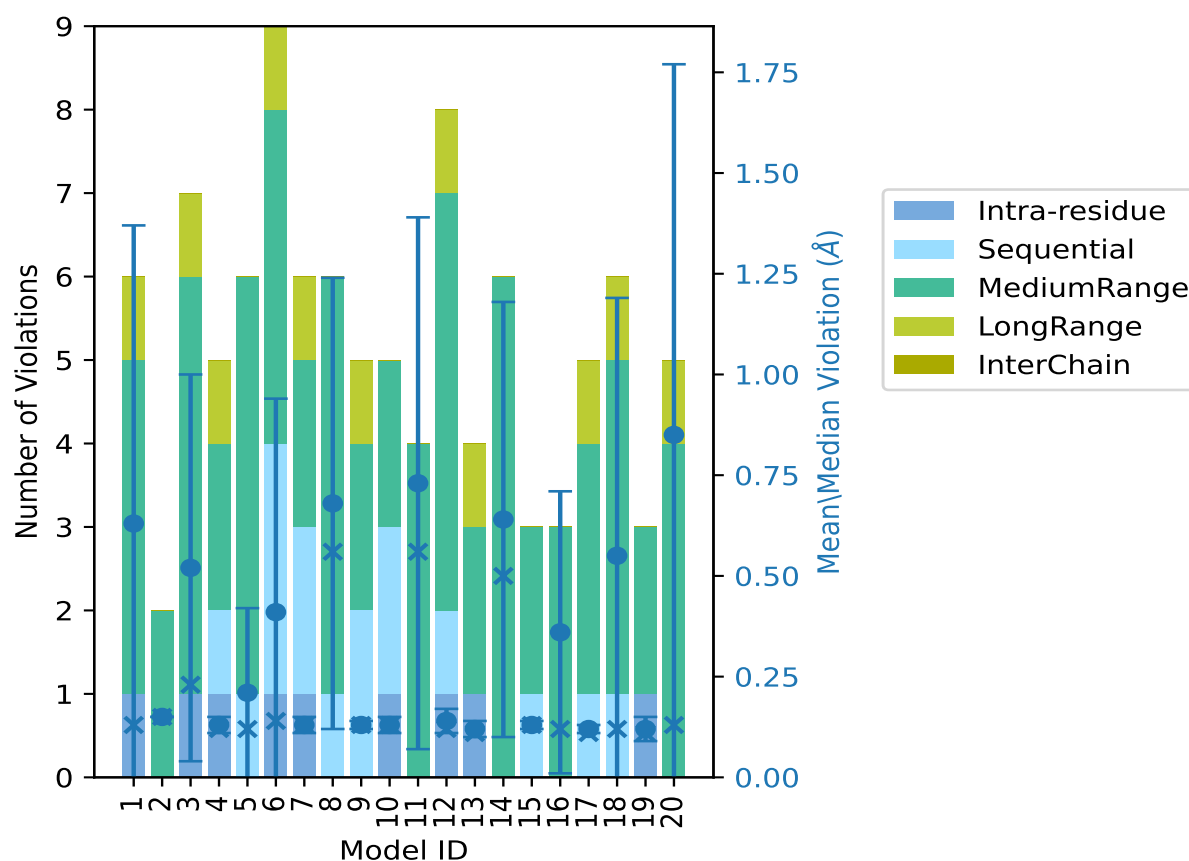
Continued from previous page...

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				
11	0	0	4	0	0	4	0.73	1.69	0.66	0.56
12	1	1	5	1	0	8	0.14	0.21	0.03	0.12
13	1	0	2	1	0	4	0.12	0.16	0.02	0.11
14	0	0	6	0	0	6	0.64	1.55	0.54	0.5
15	0	1	2	0	0	3	0.13	0.15	0.01	0.13
16	0	0	3	0	0	3	0.36	0.86	0.35	0.12
17	0	1	3	1	0	5	0.12	0.14	0.01	0.11
18	0	1	4	1	0	6	0.55	1.76	0.64	0.12
19	1	0	2	0	0	3	0.12	0.16	0.03	0.11
20	0	0	4	1	0	5	0.85	2.3	0.92	0.13

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

⁵Inter-chain restraints, ⁶Standard deviation

9.2.1 Bar graph : Distance Violation statistics for each model ⓘ



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

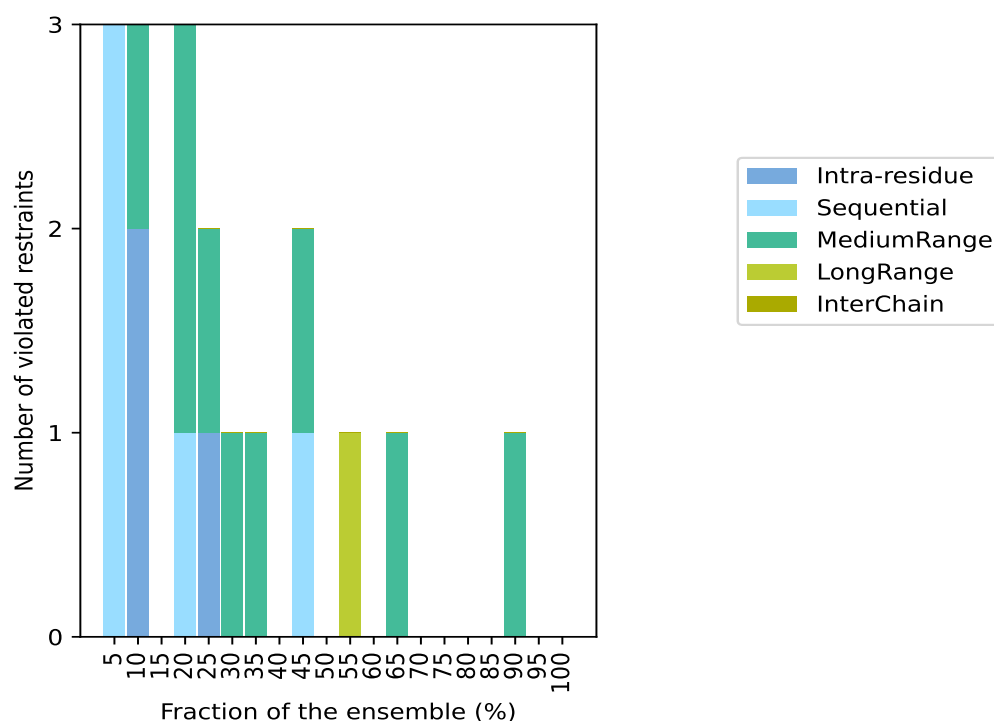
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 2524(IR:756, SQ:751, MR:299, LR:718, IC:0) restraints are 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	3	0	0	0	3	1	5.0
2	0	1	0	0	3	2	10.0
0	0	0	0	0	0	3	15.0
0	1	2	0	0	3	4	20.0
1	0	1	0	0	2	5	25.0
0	0	1	0	0	1	6	30.0
0	0	1	0	0	1	7	35.0
0	0	0	0	0	0	8	40.0
0	1	1	0	0	2	9	45.0
0	0	0	0	0	0	10	50.0
0	0	0	1	0	1	11	55.0
0	0	0	0	0	0	12	60.0
0	0	1	0	0	1	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	1	0	0	1	18	90.0
0	0	0	0	0	0	19	95.0
0	0	0	0	0	0	20	100.0

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

⁵Inter-chain restraints, ⁶ Number of models with violations

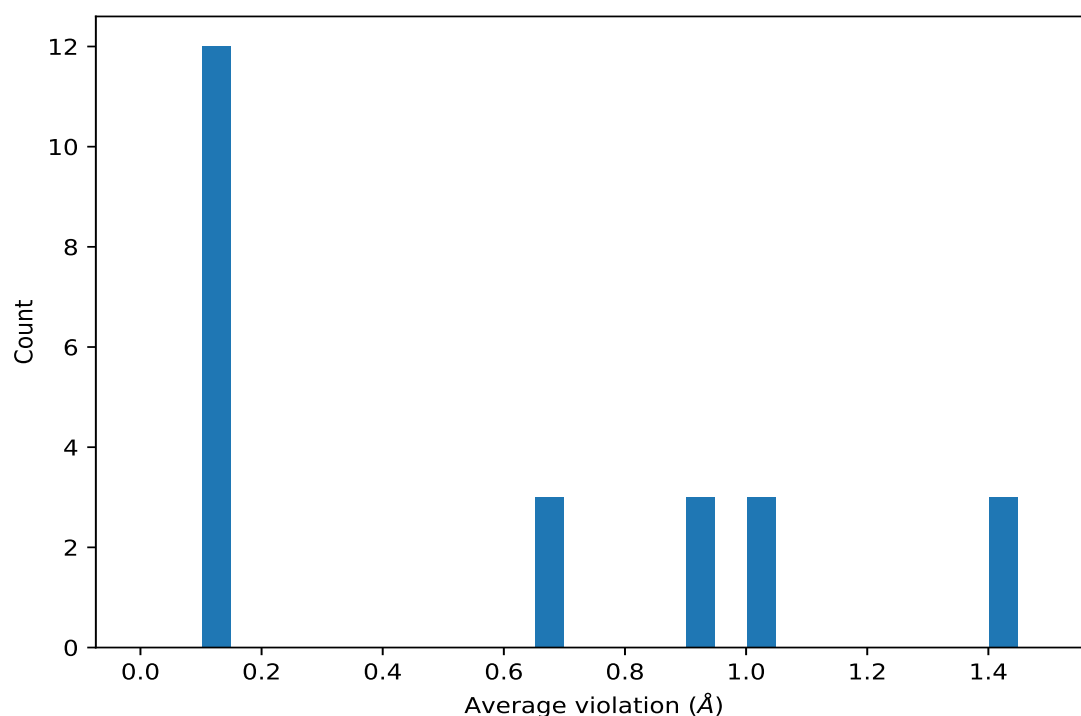
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



9.4 Most violated distance restraints in the ensemble [i](#)

9.4.1 Histogram : Distribution of mean distance violations [i](#)

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models in the ensemble



9.4.2 Table: Most violated distance restraints [i](#)

The following table provides the mean and the standard deviation of the violation for each restraint sorted by number of violated models and the mean value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	18	0.14	0.01	0.14
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	13	0.13	0.01	0.13
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	11	0.13	0.01	0.13
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	9	1.04	0.65	1.0
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	9	1.04	0.65	1.0
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	9	1.04	0.65	1.0
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	9	0.11	0.0	0.11
(4,888)	1:165:A:LEU:HD11	1:168:A:PHE:HE1	7	1.43	0.53	1.58
(4,888)	1:165:A:LEU:HD12	1:168:A:PHE:HE1	7	1.43	0.53	1.58
(4,888)	1:165:A:LEU:HD13	1:168:A:PHE:HE1	7	1.43	0.53	1.58
(4,885)	1:165:A:LEU:HD11	1:168:A:PHE:HD1	6	0.92	0.21	0.92
(4,885)	1:165:A:LEU:HD12	1:168:A:PHE:HD1	6	0.92	0.21	0.92
(4,885)	1:165:A:LEU:HD13	1:168:A:PHE:HD1	6	0.92	0.21	0.92
(4,886)	1:165:A:LEU:HD21	1:168:A:PHE:HD1	5	0.66	0.56	0.31
(4,886)	1:165:A:LEU:HD22	1:168:A:PHE:HD1	5	0.66	0.56	0.31
(4,886)	1:165:A:LEU:HD23	1:168:A:PHE:HD1	5	0.66	0.56	0.31

Continued on next page...

Continued from previous page...

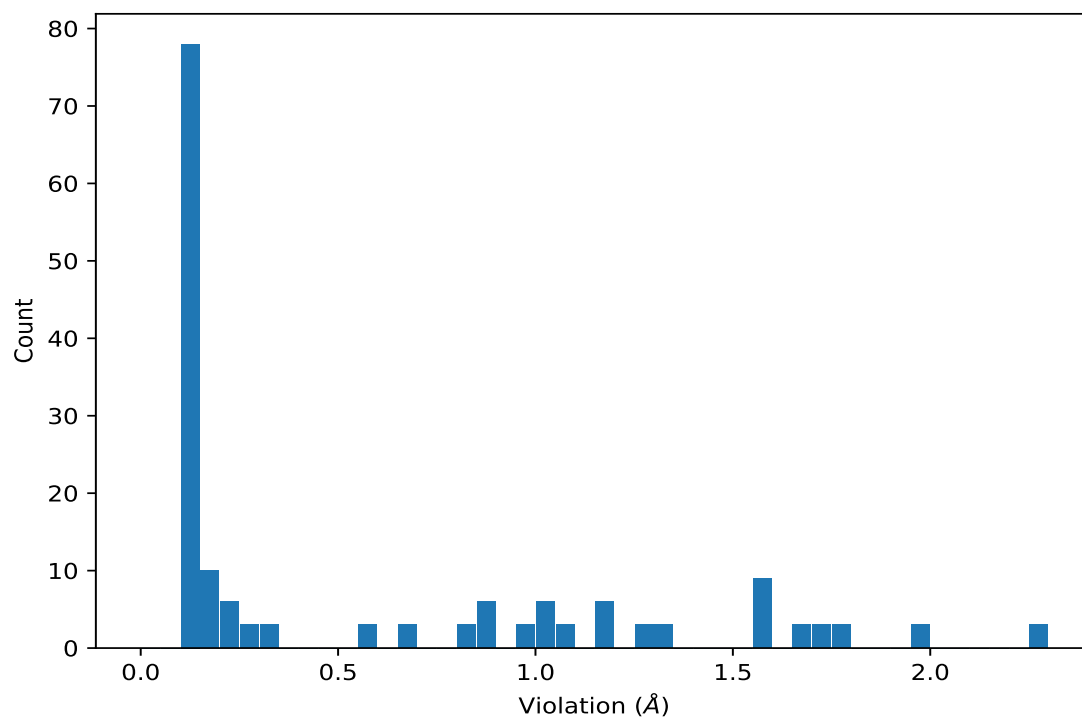
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE1	5	0.12	0.01	0.12
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE2	5	0.12	0.01	0.12
(4,948)	1:177:A:LEU:HG	1:178:A:HIS:HA	4	0.15	0.0	0.15
(3,864)	1:96:A:ARG:HA	1:93:A:LEU:H	4	0.11	0.0	0.11
(3,904)	1:97:A:GLY:H	1:99:A:CYS:H	4	0.11	0.0	0.11
(3,1473)	1:173:A:THR:HB	1:176:A:THR:H	2	0.12	0.0	0.12
(3,1532)	1:182:A:ILE:HB	1:182:A:ILE:H	2	0.11	0.0	0.11
(3,194)	1:25:A:VAL:HB	1:25:A:VAL:H	2	0.1	0.0	0.1

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)
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	20	2.3
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	20	2.3
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	20	2.3
(4,888)	1:165:A:LEU:HD11	1:168:A:PHE:HE1	1	1.99
(4,888)	1:165:A:LEU:HD12	1:168:A:PHE:HE1	1	1.99
(4,888)	1:165:A:LEU:HD13	1:168:A:PHE:HE1	1	1.99
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	18	1.76
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	18	1.76
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	18	1.76
(4,888)	1:165:A:LEU:HD11	1:168:A:PHE:HE1	6	1.71
(4,888)	1:165:A:LEU:HD12	1:168:A:PHE:HE1	6	1.71
(4,888)	1:165:A:LEU:HD13	1:168:A:PHE:HE1	6	1.71
(4,888)	1:165:A:LEU:HD11	1:168:A:PHE:HE1	11	1.69
(4,888)	1:165:A:LEU:HD12	1:168:A:PHE:HE1	11	1.69
(4,888)	1:165:A:LEU:HD13	1:168:A:PHE:HE1	11	1.69
(4,888)	1:165:A:LEU:HD11	1:168:A:PHE:HE1	8	1.58
(4,888)	1:165:A:LEU:HD12	1:168:A:PHE:HE1	8	1.58
(4,888)	1:165:A:LEU:HD13	1:168:A:PHE:HE1	8	1.58
(4,886)	1:165:A:LEU:HD21	1:168:A:PHE:HD1	20	1.56
(4,886)	1:165:A:LEU:HD22	1:168:A:PHE:HD1	20	1.56
(4,886)	1:165:A:LEU:HD23	1:168:A:PHE:HD1	20	1.56
(4,888)	1:165:A:LEU:HD11	1:168:A:PHE:HE1	14	1.55
(4,888)	1:165:A:LEU:HD12	1:168:A:PHE:HE1	14	1.55
(4,888)	1:165:A:LEU:HD13	1:168:A:PHE:HE1	14	1.55
(4,888)	1:165:A:LEU:HD11	1:168:A:PHE:HE1	3	1.31
(4,888)	1:165:A:LEU:HD12	1:168:A:PHE:HE1	3	1.31
(4,888)	1:165:A:LEU:HD13	1:168:A:PHE:HE1	3	1.31
(4,885)	1:165:A:LEU:HD11	1:168:A:PHE:HD1	1	1.27
(4,885)	1:165:A:LEU:HD12	1:168:A:PHE:HD1	1	1.27
(4,885)	1:165:A:LEU:HD13	1:168:A:PHE:HD1	1	1.27
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	8	1.19
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	8	1.19
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	8	1.19
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	3	1.18
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	3	1.18
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	3	1.18
(4,886)	1:165:A:LEU:HD21	1:168:A:PHE:HD1	18	1.07
(4,886)	1:165:A:LEU:HD22	1:168:A:PHE:HD1	18	1.07
(4,886)	1:165:A:LEU:HD23	1:168:A:PHE:HD1	18	1.07
(4,885)	1:165:A:LEU:HD11	1:168:A:PHE:HD1	6	1.01
(4,885)	1:165:A:LEU:HD12	1:168:A:PHE:HD1	6	1.01

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(4,885)	1:165:A:LEU:HD13	1:168:A:PHE:HD1	6	1.01
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	14	1.0
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	14	1.0
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	14	1.0
(4,885)	1:165:A:LEU:HD11	1:168:A:PHE:HD1	11	0.99
(4,885)	1:165:A:LEU:HD12	1:168:A:PHE:HD1	11	0.99
(4,885)	1:165:A:LEU:HD13	1:168:A:PHE:HD1	11	0.99
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	16	0.86
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	16	0.86
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	16	0.86
(4,885)	1:165:A:LEU:HD11	1:168:A:PHE:HD1	14	0.85
(4,885)	1:165:A:LEU:HD12	1:168:A:PHE:HD1	14	0.85
(4,885)	1:165:A:LEU:HD13	1:168:A:PHE:HD1	14	0.85
(4,885)	1:165:A:LEU:HD11	1:168:A:PHE:HD1	8	0.82
(4,885)	1:165:A:LEU:HD12	1:168:A:PHE:HD1	8	0.82
(4,885)	1:165:A:LEU:HD13	1:168:A:PHE:HD1	8	0.82
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	5	0.67
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	5	0.67
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	5	0.67
(4,885)	1:165:A:LEU:HD11	1:168:A:PHE:HD1	3	0.58
(4,885)	1:165:A:LEU:HD12	1:168:A:PHE:HD1	3	0.58
(4,885)	1:165:A:LEU:HD13	1:168:A:PHE:HD1	3	0.58
(4,886)	1:165:A:LEU:HD21	1:168:A:PHE:HD1	8	0.31
(4,886)	1:165:A:LEU:HD22	1:168:A:PHE:HD1	8	0.31
(4,886)	1:165:A:LEU:HD23	1:168:A:PHE:HD1	8	0.31
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	6	0.26
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	6	0.26
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	6	0.26
(4,886)	1:165:A:LEU:HD21	1:168:A:PHE:HD1	3	0.23
(4,886)	1:165:A:LEU:HD22	1:168:A:PHE:HD1	3	0.23
(4,886)	1:165:A:LEU:HD23	1:168:A:PHE:HD1	3	0.23
(4,888)	1:165:A:LEU:HD11	1:168:A:PHE:HE1	12	0.21
(4,888)	1:165:A:LEU:HD12	1:168:A:PHE:HE1	12	0.21
(4,888)	1:165:A:LEU:HD13	1:168:A:PHE:HE1	12	0.21
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	13	0.16
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	19	0.16
(4,948)	1:177:A:LEU:HG	1:178:A:HIS:HA	10	0.15
(4,948)	1:177:A:LEU:HG	1:178:A:HIS:HA	15	0.15
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	2	0.15
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	4	0.15
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	7	0.15
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	9	0.15

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	10	0.15
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	14	0.15
(4,948)	1:177:A:LEU:HG	1:178:A:HIS:HA	6	0.14
(4,948)	1:177:A:LEU:HG	1:178:A:HIS:HA	9	0.14
(4,886)	1:165:A:LEU:HD21	1:168:A:PHE:HD1	14	0.14
(4,886)	1:165:A:LEU:HD22	1:168:A:PHE:HD1	14	0.14
(4,886)	1:165:A:LEU:HD23	1:168:A:PHE:HD1	14	0.14
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	4	0.14
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	7	0.14
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	12	0.14
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	2	0.14
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	7	0.14
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	14	0.14
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	17	0.14
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	6	0.14
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	12	0.14
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE1	1	0.13
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE2	1	0.13
(3,1080)	1:118:A:VAL:HB	1:119:A:ALA:H	6	0.13
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	3	0.13
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	9	0.13
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	5	0.13
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	9	0.13
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	10	0.13
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	12	0.13
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	1	0.13
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	5	0.13
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	11	0.13
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	15	0.13
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	20	0.13
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE1	10	0.12
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE2	10	0.12
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE1	12	0.12
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE2	12	0.12
(3,1473)	1:173:A:THR:HB	1:176:A:THR:H	5	0.12
(3,1473)	1:173:A:THR:HB	1:176:A:THR:H	15	0.12
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	4	0.12
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	7	0.12
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	12	0.12
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	1	0.12
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	6	0.12
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	17	0.12

Continued on next page...

Continued from previous page...

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	18	0.12
(3,1020)	1:78:A:ARG:HA	1:110:A:GLN:H	20	0.12
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	1	0.12
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	4	0.12
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	11	0.12
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	18	0.12
(3,903)	1:96:A:ARG:HA	1:99:A:CYS:H	20	0.12
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	16	0.12
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	18	0.12
(4,889)	1:165:A:LEU:HD21	1:168:A:PHE:HE1	12	0.11
(4,889)	1:165:A:LEU:HD22	1:168:A:PHE:HE1	12	0.11
(4,889)	1:165:A:LEU:HD23	1:168:A:PHE:HE1	12	0.11
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE1	13	0.11
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE2	13	0.11
(3,1532)	1:182:A:ILE:HB	1:182:A:ILE:H	6	0.11
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	5	0.11
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	6	0.11
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	8	0.11
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	9	0.11
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	17	0.11
(3,1038)	1:111:A:ILE:HB	1:112:A:ALA:H	18	0.11
(3,904)	1:97:A:GLY:H	1:99:A:CYS:H	3	0.11
(3,904)	1:97:A:GLY:H	1:99:A:CYS:H	13	0.11
(3,904)	1:97:A:GLY:H	1:99:A:CYS:H	16	0.11
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	13	0.11
(3,873)	1:94:A:ALA:H	1:96:A:ARG:HA	17	0.11
(3,864)	1:96:A:ARG:HA	1:93:A:LEU:H	5	0.11
(3,864)	1:96:A:ARG:HA	1:93:A:LEU:H	12	0.11
(3,864)	1:96:A:ARG:HA	1:93:A:LEU:H	17	0.11
(3,864)	1:96:A:ARG:HA	1:93:A:LEU:H	19	0.11
(3,305)	1:35:A:GLN:HB2	1:36:A:GLU:H	10	0.11
(3,304)	1:35:A:GLN:HB3	1:36:A:GLU:H	7	0.11
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE1	19	0.1
(4,791)	1:134:A:TYR:HA	1:134:A:TYR:HE2	19	0.1
(3,1532)	1:182:A:ILE:HB	1:182:A:ILE:H	4	0.1
(3,904)	1:97:A:GLY:H	1:99:A:CYS:H	8	0.1
(3,194)	1:25:A:VAL:HB	1:25:A:VAL:H	3	0.1
(3,194)	1:25:A:VAL:HB	1:25:A:VAL:H	7	0.1

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

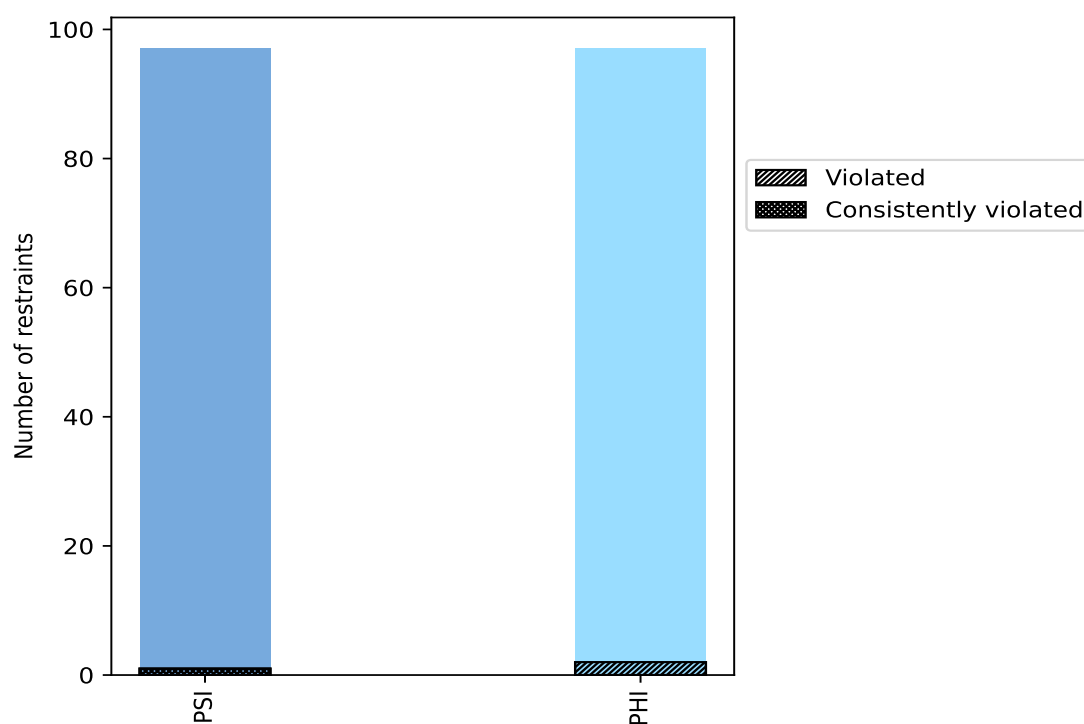
10.1 Summary of dihedral-angle violations [i](#)

The following table provides the summary of dihedral-angle violations in different dihedral-angle types. Violations less than 1° are not included in the calculation.

Angle type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
PSI	97	50.0	1	1.0	0.5	1	1.0	0.5
PHI	97	50.0	2	2.1	1.0	0	0.0	0.0
Total	194	100.0	3	1.5	1.5	1	0.5	0.5

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](#)



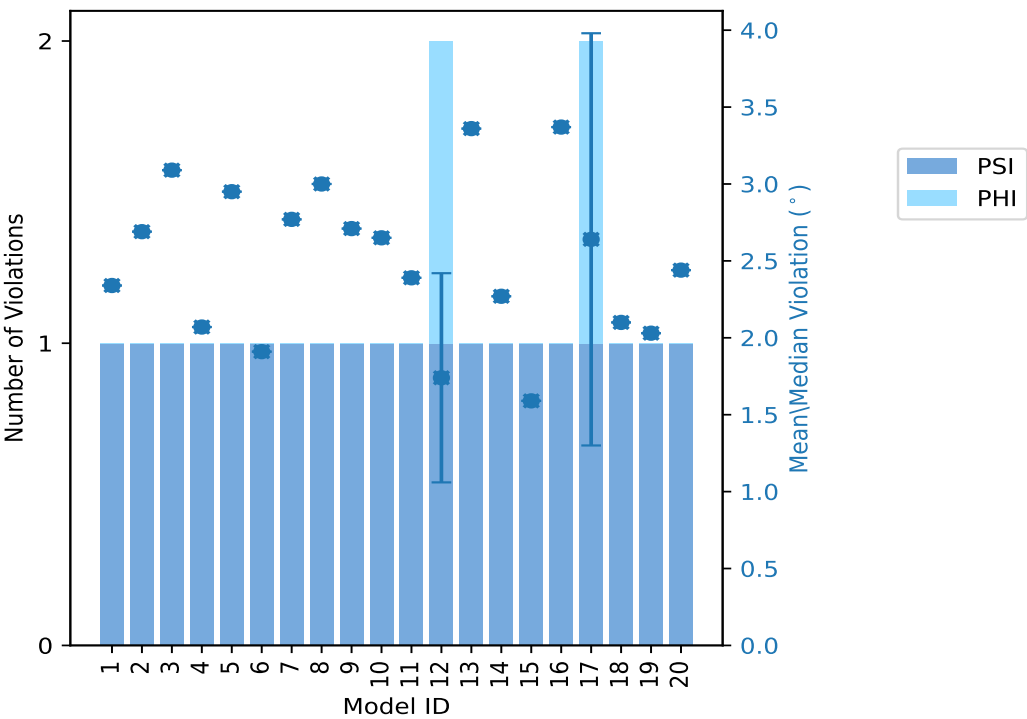
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

10.2 Dihedral-angle violation statistics for each model

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	1	0	1	2.34	2.34	0.0	2.34
2	1	0	1	2.69	2.69	0.0	2.69
3	1	0	1	3.09	3.09	0.0	3.09
4	1	0	1	2.07	2.07	0.0	2.07
5	1	0	1	2.95	2.95	0.0	2.95
6	1	0	1	1.91	1.91	0.0	1.91
7	1	0	1	2.77	2.77	0.0	2.77
8	1	0	1	3.0	3.0	0.0	3.0
9	1	0	1	2.71	2.71	0.0	2.71
10	1	0	1	2.65	2.65	0.0	2.65
11	1	0	1	2.39	2.39	0.0	2.39
12	1	1	2	1.74	2.41	0.68	1.74
13	1	0	1	3.36	3.36	0.0	3.36
14	1	0	1	2.27	2.27	0.0	2.27
15	1	0	1	1.59	1.59	0.0	1.59
16	1	0	1	3.37	3.37	0.0	3.37
17	1	1	2	2.64	3.97	1.34	2.64
18	1	0	1	2.1	2.1	0.0	2.1
19	1	0	1	2.03	2.03	0.0	2.03
20	1	0	1	2.44	2.44	0.0	2.44

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	
PSI	PHI	Total	Count ¹	%
0	2	2	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

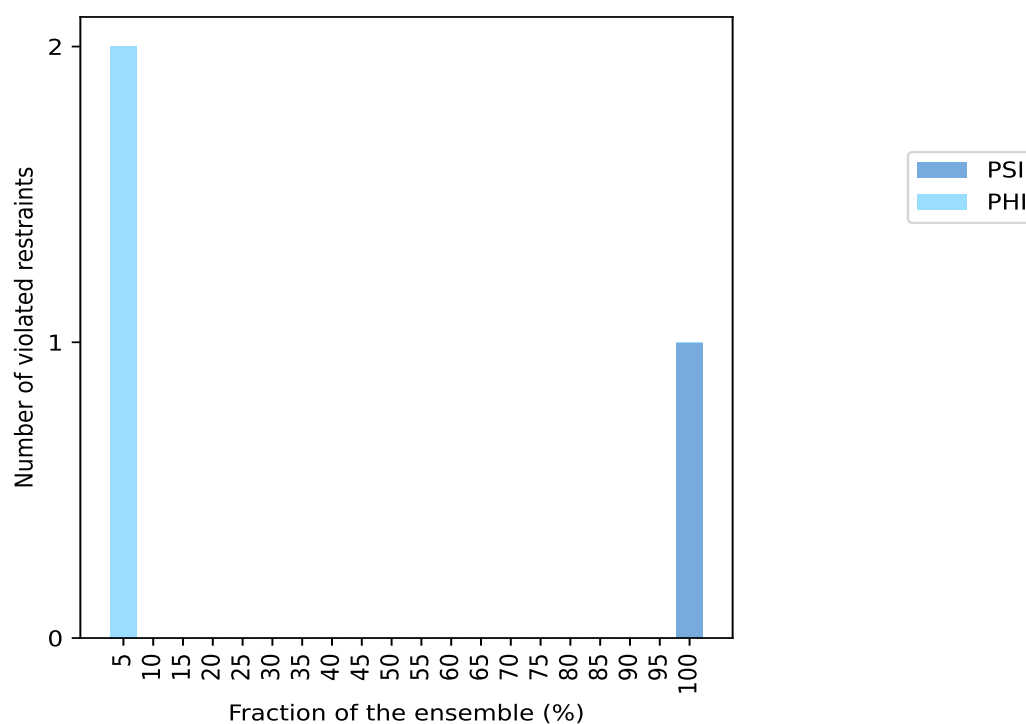
Continued on next page...

Continued from previous page...

Number of violated restraints			Fraction of the ensemble	
PSI	PHI	Total	Count ¹	%
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
1	0	1	20	100.0

¹ Number of models with violations

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

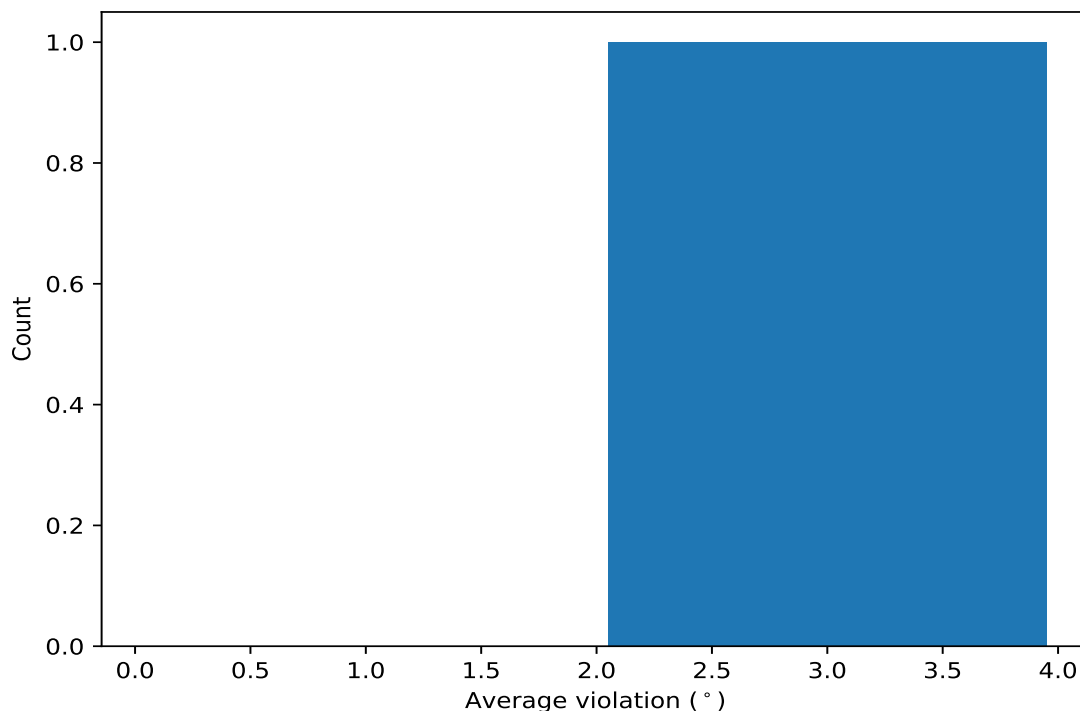


10.4 Most violated dihedral-angle restraints in the ensemble [i](#)

10.4.1 Histogram : Distribution of mean dihedral-angle violations [i](#)

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models

in the ensemble



10.4.2 Table: Most violated dihedral-angle restraints [i](#)

The following table provides the mean and the standard deviation of the 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.

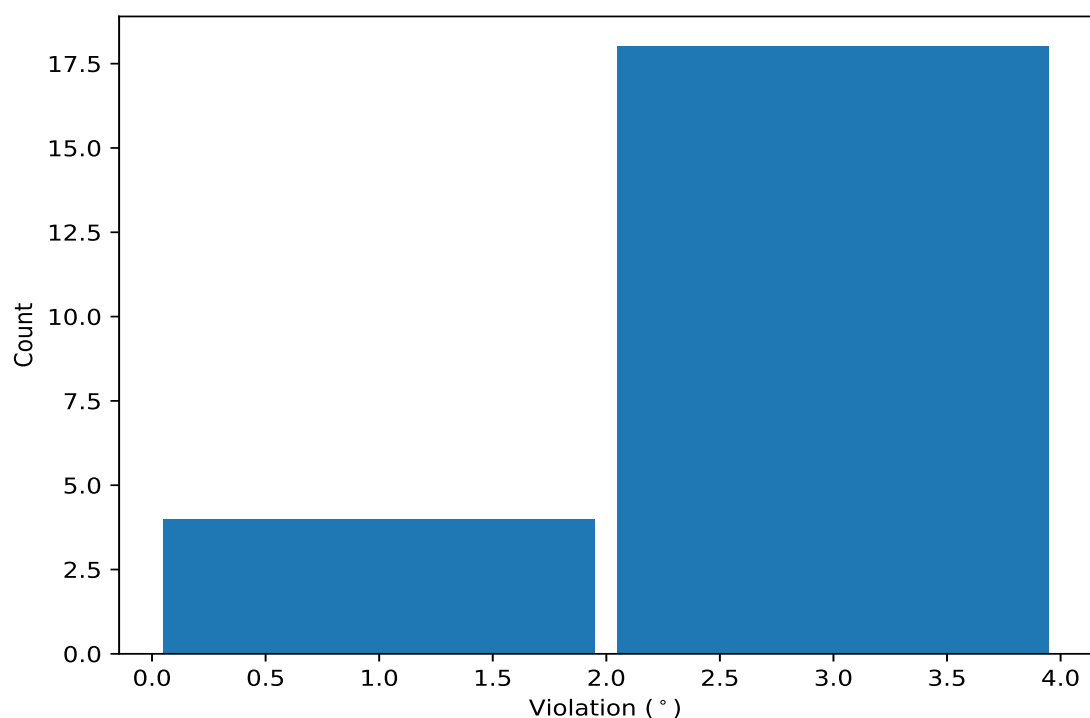
Key	Atom-1	Atom-2	Atom-3	Atom-4	Models ¹	Mean	SD ²	Median
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	20	2.61	0.56	2.54

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

10.5 All violated dihedral-angle restraints [i](#)

10.5.1 Histogram : Distribution of violations [i](#)

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



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

The following table 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 (°)
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	17	3.97
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	16	3.37
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	13	3.36
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	3	3.09
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	8	3.0
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	5	2.95
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	7	2.77
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	9	2.71
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	2	2.69
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	10	2.65
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	20	2.44
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	12	2.41
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	11	2.39
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	1	2.34
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	14	2.27
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	18	2.1
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	4	2.07
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	19	2.03
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	6	1.91
(1,108)	1:98:A:ARG:N	1:98:A:ARG:CA	1:98:A:ARG:C	1:99:A:CYS:N	15	1.59
(1,109)	1:98:A:ARG:C	1:99:A:CYS:N	1:99:A:CYS:CA	1:99:A:CYS:C	17	1.3
(1,129)	1:110:A:GLN:C	1:111:A:ILE:N	1:111:A:ILE:CA	1:111:A:ILE:C	12	1.06