



wwPDB NMR Structure Validation Summary Report ⓘ

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BMRB ID : 35007
Title : Hybrid structure of the helix bundle domain of pRN1 primase in complex with DNA, ATP and dGTP
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<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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
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) : 2.49

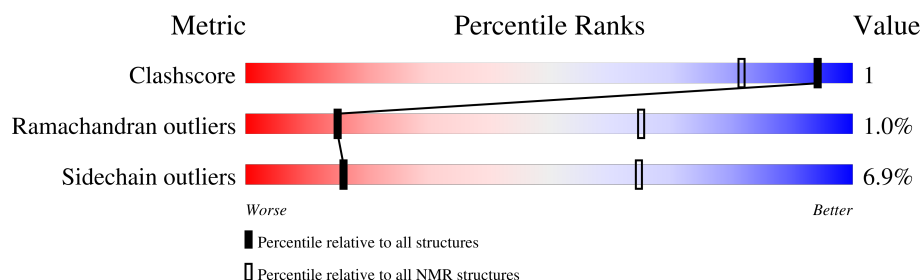
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR, SOLID-STATE NMR

The overall completeness of chemical shifts assignment is 89%.

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	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

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	B	118	
2	A	9	

2 Ensemble composition and analysis

This entry contains 20 models. Model 14 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	B:260-B:370 (111)	0.18	14

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

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

The models can be grouped into 3 clusters and 10 single-model clusters were found.

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

3 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2345 atoms, of which 1130 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called SF3 helicase domain-containing protein.

Mol	Chain	Residues	Atoms						Trace
1	B	118	Total	C	H	N	O	S	0
			1978	624	1006	165	179	4	

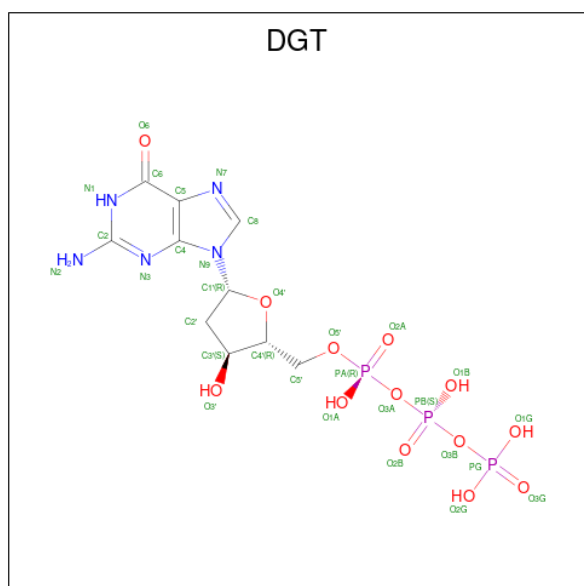
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	253	GLY	-	expression tag	UNP Q54324
B	254	HIS	-	expression tag	UNP Q54324
B	255	MET	-	expression tag	UNP Q54324

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*TP*GP*TP*GP*CP*TP*CP*A)-3').

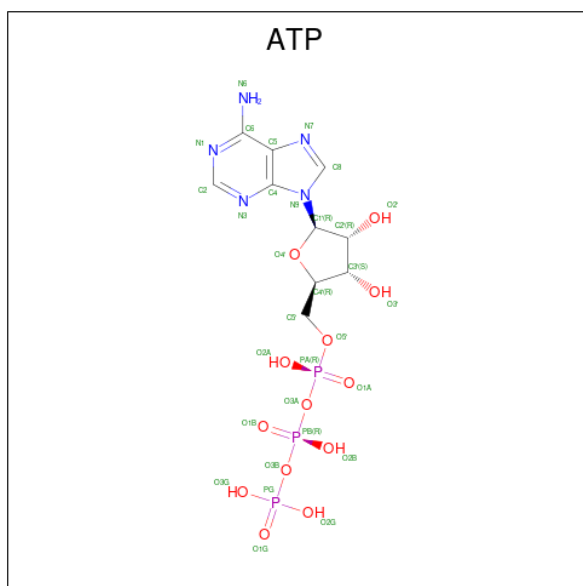
Mol	Chain	Residues	Atoms						Trace
2	A	9	Total	C	H	N	O	P	0
			281	87	102	30	54	8	

- Molecule 3 is 2'-DEOXYGUANOSINE-5'-TRIPHOSPHATE (CCD ID: DGT) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					
3	B	1	Total	C	H	N	O	P
			42	10	11	5	13	3

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					
4	B	1	Total	C	H	N	O	P
			42	10	11	5	13	3

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

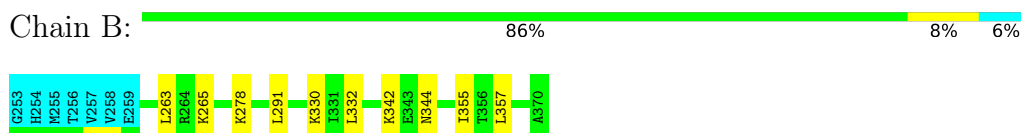
Mol	Chain	Residues	Atoms	
5	B	2	Total	Mg
			2	2

4 Residue-property plots [i](#)

4.1 Average score per residue in the NMR ensemble

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

- Molecule 1: SF3 helicase domain-containing protein



- Molecule 2: DNA (5'-D(*CP*TP*GP*TP*GP*CP*TP*CP*A)-3')

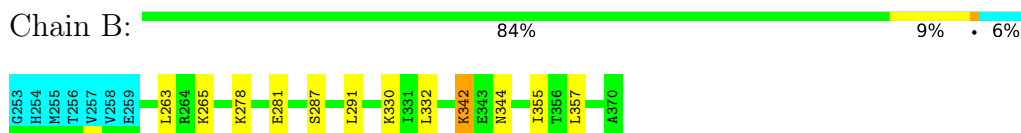


There are no outlier residues in this chain.

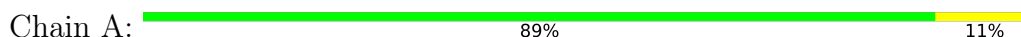
4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 14. Colouring as in section 4.1 above.

- Molecule 1: SF3 helicase domain-containing protein



- Molecule 2: DNA (5'-D(*CP*TP*GP*TP*GP*CP*TP*CP*A)-3')



5 Refinement protocol and experimental data overview

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

Of the 500 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
CYANA	structure calculation	3.98.15
Amber	refinement	20

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	4
Total number of shifts	1694
Number of shifts mapped to atoms	1694
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	89%

Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

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: ATP, MG, DGT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	B	0.52±0.00	0±0/938 (0.0± 0.0%)	0.96±0.02	1±1/1261 (0.0± 0.1%)
2	A	0.51±0.01	0±0/199 (0.0± 0.0%)	1.03±0.03	0±0/305 (0.0± 0.1%)
All	All	0.52	0/22740 (0.0%)	0.97	13/31320 (0.0%)

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	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	B	287	SER	CA-C-N	5.37	123.53	119.66	9	6
1	B	287	SER	C-N-CA	5.37	123.53	119.66	9	6
2	A	4	DT	C5'-C4'-C3'	-5.06	107.31	114.90	8	1

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	B	920	956	956	2±1
2	A	179	102	104	0±1
4	B	31	11	12	0±0
All	All	23260	21600	21678	54

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 1.

5 of 12 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:340:LYS:HE3	1:B:346:LYS:HE2	0.49	1.85	19	1
1:B:330:LYS:HE2	1:B:330:LYS:HA	0.49	1.84	19	20
1:B:342:LYS:HB2	1:B:342:LYS:HZ2	0.44	1.72	11	3
1:B:291:LEU:HD13	1:B:291:LEU:C	0.43	2.38	3	13
1:B:340:LYS:CE	1:B:346:LYS:HE2	0.43	2.43	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	B	110/118 (93%)	104±1 (95±1%)	5±1 (4±1%)	1±0 (1±0%)	15	65
All	All	2200/2360 (93%)	2081 (95%)	96 (4%)	23 (1%)	15	65

All 2 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	B	344	ASN	20
1	B	339	SER	3

6.3.2 Protein sidechains [i](#)

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	104/110 (95%)	97±1 (93±1%)	7±1 (7±1%)	16	65
All	All	2080/2200 (95%)	1936 (93%)	144 (7%)	16	65

5 of 9 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	B	263	LEU	20
1	B	265	LYS	20
1	B	278	LYS	20
1	B	332	LEU	20
1	B	342	LYS	20

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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
3	DGT	B	401	5	32,33,33	0.74±0.01	0±0 (0±0%)
4	ATP	B	402	5	32,33,33	0.76±0.01	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.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
3	DGT	B	401	5	48,52,52	0.76±0.01	0±0 (0±0%)
4	ATP	B	402	5	48,52,52	0.78±0.03	1±0 (1±0%)

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
3	DGT	B	401	5	-	0±0,22,34,34	0±0,3,3,3
4	ATP	B	402	5	-	0±0,22,38,38	0±0,3,3,3

There are no bond-length outliers.

All unique angle outliers are listed below.

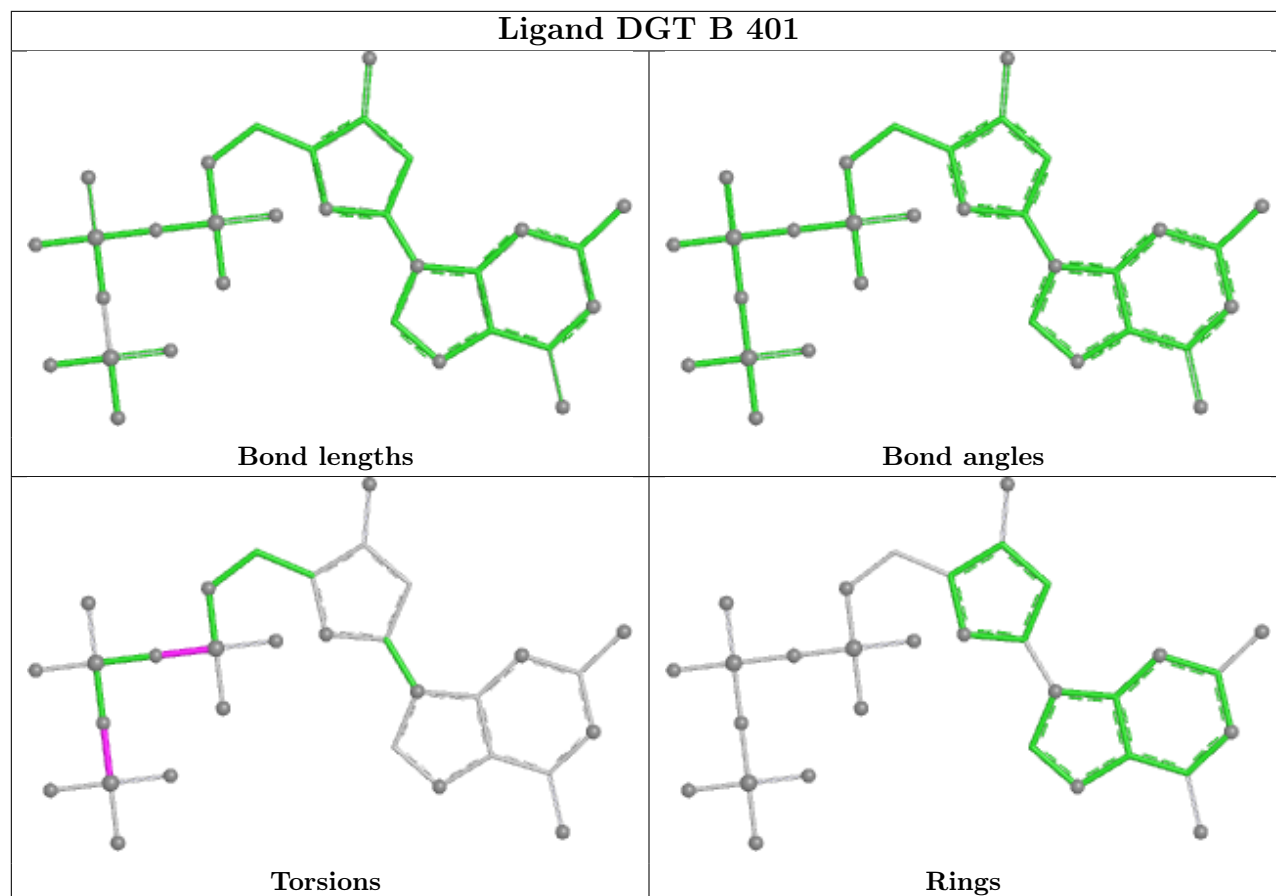
Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
4	B	402	ATP	O3G-PG-O2G	2.09	115.64	107.80	12	15

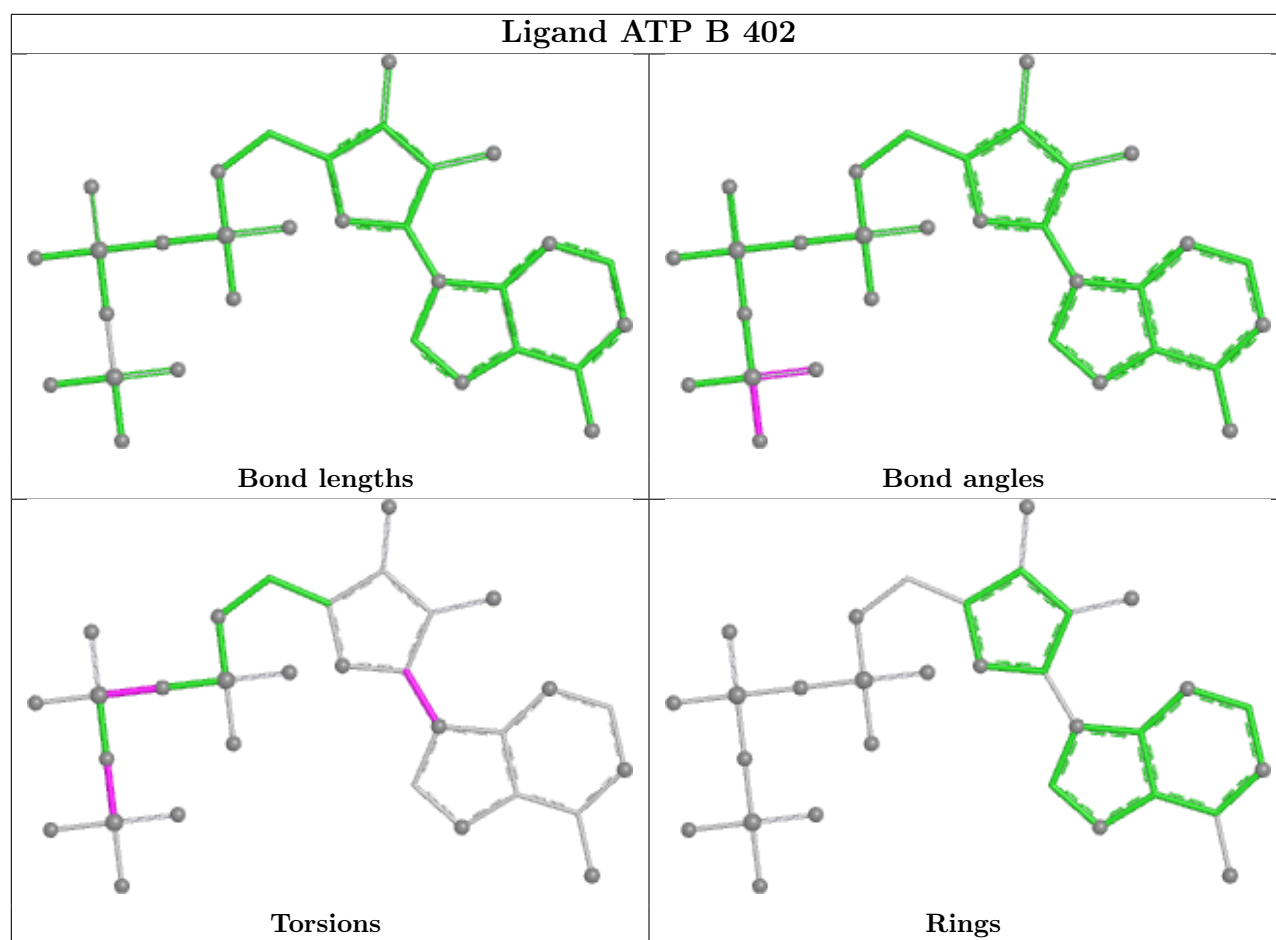
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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 89% for the well-defined parts and 89% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chemical_shifts_DNA*

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	101
Number of shifts mapped to atoms	101
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	3

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 5%, i.e. 92 atoms were assigned a chemical shift out of a possible 1789. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/548 (0%)	0/220 (0%)	0/222 (0%)	0/106 (0%)
Sidechain	0/959 (0%)	0/617 (0%)	0/304 (0%)	0/38 (0%)
Aromatic	0/106 (0%)	0/52 (0%)	0/49 (0%)	0/5 (0%)
Sugar	71/108 (66%)	63/63 (100%)	8/45 (18%)	0/0 (—%)
Base	21/68 (31%)	16/41 (39%)	2/16 (12%)	3/11 (27%)
Overall	92/1789 (5%)	79/993 (8%)	10/636 (2%)	3/160 (2%)

Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a

chemical shift value, which may lead to lower completeness of assignment measure.

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	3	DG	H8	5.50	6.49 – 9.09	-8.8
1	A	5	DG	H2''	1.17	1.57 – 3.80	-6.8
1	A	4	DT	H2'	0.58	0.76 – 3.44	-5.7

7.1.5 Random Coil Index (RCI) plots [i](#)

No *random coil index*(RCI) plot could be generated from the current chemical shift list. RCI is only applicable to proteins

7.2 Chemical shift list 2

File name: working_cs.cif

Chemical shift list name: *assigned_chemical_shifts_protein*

7.2.1 Bookkeeping [i](#)

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

Total number of shifts	1581
Number of shifts mapped to atoms	1581
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	4

7.2.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	117	-0.53 \pm 0.22	Should be checked

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Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\beta$	112	0.43 ± 0.05	None needed (< 0.5 ppm)
$^{13}\text{C}'$	114	-0.16 ± 0.21	None needed (< 0.5 ppm)
^{15}N	108	1.05 ± 0.31	Should be applied

7.2.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 84%, i.e. 1498 atoms were assigned a chemical shift out of a possible 1789. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	539/548 (98%)	218/220 (99%)	218/222 (98%)	103/106 (97%)
Sidechain	876/959 (91%)	593/617 (96%)	273/304 (90%)	10/38 (26%)
Aromatic	83/106 (78%)	46/52 (88%)	34/49 (69%)	3/5 (60%)
Sugar	0/108 (0%)	0/63 (0%)	0/45 (0%)	0/0 (—%)
Base	0/68 (0%)	0/41 (0%)	0/16 (0%)	0/11 (0%)
Overall	1498/1789 (84%)	857/993 (86%)	525/636 (83%)	116/160 (72%)

Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

7.2.4 Statistically unusual chemical shifts [i](#)

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

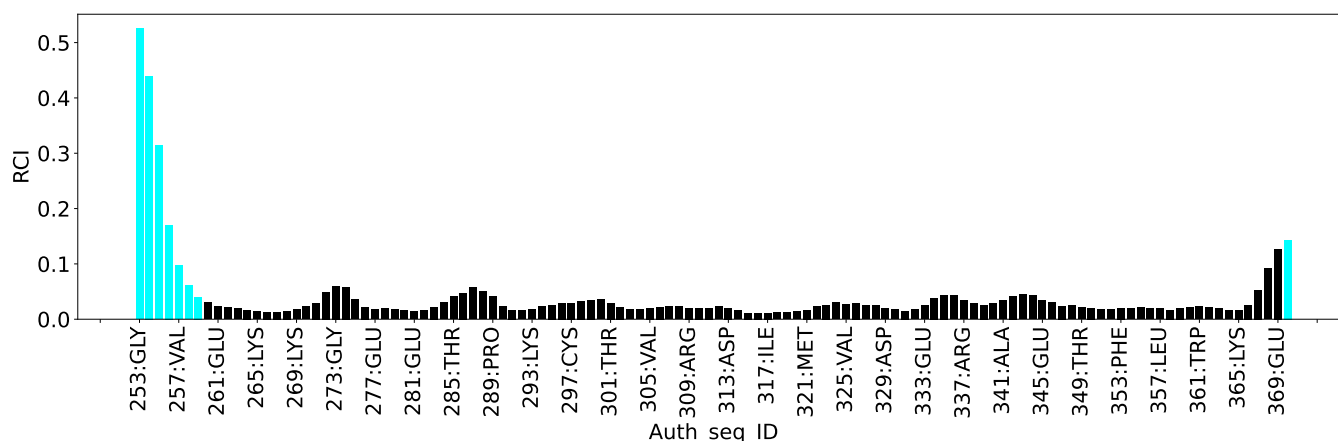
List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
2	B	356	THR	HG1	5.49	0.08 – 2.19	20.6
2	B	253	GLY	CA	63.91	38.93 – 51.79	14.4
2	B	351	LYS	HB2	0.30	0.58 – 2.97	-6.2
2	B	351	LYS	HB3	0.40	0.46 – 3.04	-5.2

7.2.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble

composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain B:



7.3 Chemical shift list 3

File name: working_cs.cif

Chemical shift list name: *chem_shift_list_DGT*

7.3.1 Bookkeeping [i](#)

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

Total number of shifts	7
Number of shifts mapped to atoms	7
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

7.3.2 Chemical shift referencing [i](#)

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

7.3.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 0%, i.e. 0 atoms were assigned a chemical

shift out of a possible 1789. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	0/548 (0%)	0/220 (0%)	0/222 (0%)	0/106 (0%)
Sidechain	0/959 (0%)	0/617 (0%)	0/304 (0%)	0/38 (0%)
Aromatic	0/106 (0%)	0/52 (0%)	0/49 (0%)	0/5 (0%)
Sugar	0/108 (0%)	0/63 (0%)	0/45 (0%)	0/0 (—%)
Base	0/68 (0%)	0/41 (0%)	0/16 (0%)	0/11 (0%)
Overall	0/1789 (0%)	0/993 (0%)	0/636 (0%)	0/160 (0%)

Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

7.3.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.3.5 Random Coil Index (RCI) plots [i](#)

No *random coil index*(RCI) plot could be generated from the current chemical shift list. RCI is only applicable to proteins

7.4 Chemical shift list 4

File name: working_cs.cif

Chemical shift list name: *chem_shift_list_ATP*

7.4.1 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	5
Number of shifts mapped to atoms	5
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

7.4.2 Chemical shift referencing [i](#)

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

7.4.3 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 0%, i.e. 0 atoms were assigned a chemical shift out of a possible 1789. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/548 (0%)	0/220 (0%)	0/222 (0%)	0/106 (0%)
Sidechain	0/959 (0%)	0/617 (0%)	0/304 (0%)	0/38 (0%)
Aromatic	0/106 (0%)	0/52 (0%)	0/49 (0%)	0/5 (0%)
Sugar	0/108 (0%)	0/63 (0%)	0/45 (0%)	0/0 (—%)
Base	0/68 (0%)	0/41 (0%)	0/16 (0%)	0/11 (0%)
Overall	0/1789 (0%)	0/993 (0%)	0/636 (0%)	0/160 (0%)

Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

7.4.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.4.5 Random Coil Index (RCI) plots [i](#)

No *random coil index*(RCI) plot could be generated from the current chemical shift list. RCI is only applicable to proteins

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	4142
Intra-residue ($ i-j =0$)	962
Sequential ($ i-j =1$)	843
Medium range ($ i-j >1$ and $ i-j <5$)	1082
Long range ($ i-j \geq 5$)	963
Inter-chain	180
Hydrogen bond restraints	106
Disulfide bond restraints	0
Total dihedral-angle restraints	198
Number of unmapped restraints	0
Number of restraints per residue	33.1
Number of long range restraints per residue ¹	7.4

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)	63.0	0.2
0.2-0.5 (Medium)	15.4	0.5
>0.5 (Large)	8.0	1.24

8.2.2 Average number of dihedral-angle violations per model [i](#)

Dihedral-angle violations less than 1° are not included in the calculation.

Bins (°)	Average number of violations per model	Max (°)
1.0-10.0 (Small)	0.1	2.35
10.0-20.0 (Medium)	None	None
>20.0 (Large)	None	None

9 Distance violation analysis ⓘ

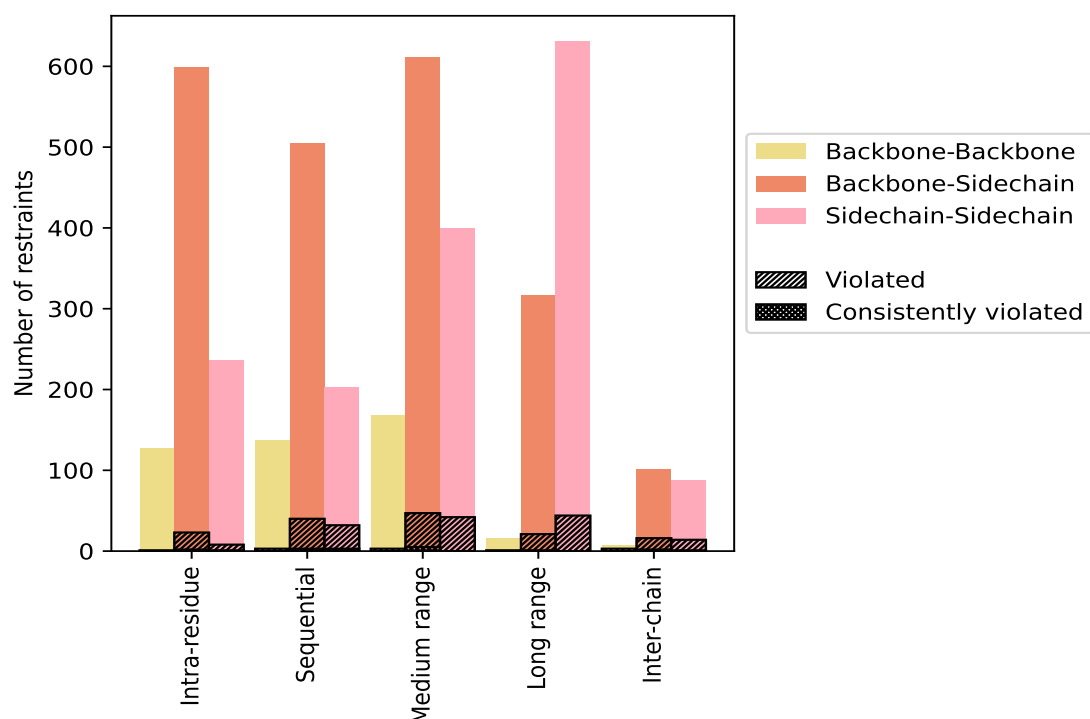
9.1 Summary of distance violations ⓘ

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)	962	23.2	32	3.3	0.8	3	0.3	0.1
Backbone-Backbone	127	3.1	1	0.8	0.0	0	0.0	0.0
Backbone-Sidechain	599	14.5	23	3.8	0.6	2	0.3	0.0
Sidechain-Sidechain	236	5.7	8	3.4	0.2	1	0.4	0.0
Sequential (i-j =1)	843	20.4	75	8.9	1.8	6	0.7	0.1
Backbone-Backbone	137	3.3	3	2.2	0.1	0	0.0	0.0
Backbone-Sidechain	504	12.2	40	7.9	1.0	3	0.6	0.1
Sidechain-Sidechain	202	4.9	32	15.8	0.8	3	1.5	0.1
Medium range (i-j >1 & i-j <5)	1082	26.1	79	7.3	1.9	0	0.0	0.0
Backbone-Backbone	168	4.1	3	1.8	0.1	0	0.0	0.0
Backbone-Sidechain	515	12.4	34	6.6	0.8	0	0.0	0.0
Sidechain-Sidechain	399	9.6	42	10.5	1.0	0	0.0	0.0
Long range (i-j ≥5)	963	23.2	66	6.9	1.6	1	0.1	0.0
Backbone-Backbone	16	0.4	1	6.2	0.0	0	0.0	0.0
Backbone-Sidechain	316	7.6	21	6.6	0.5	1	0.3	0.0
Sidechain-Sidechain	631	15.2	44	7.0	1.1	0	0.0	0.0
Inter-chain	180	4.3	29	16.1	0.7	3	1.7	0.1
Backbone-Backbone	7	0.2	3	42.9	0.1	0	0.0	0.0
Backbone-Sidechain	93	2.2	14	15.1	0.3	2	2.2	0.0
Sidechain-Sidechain	80	1.9	12	15.0	0.3	1	1.2	0.0
Hydrogen bond	106	2.6	16	15.1	0.4	5	4.7	0.1
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	4142	100.0	298	7.2	7.2	18	0.4	0.4
Backbone-Backbone	455	11.0	11	2.4	0.3	0	0.0	0.0
Backbone-Sidechain	2131	51.4	147	6.9	3.5	13	0.6	0.3
Sidechain-Sidechain	1556	37.6	140	9.0	3.4	5	0.3	0.1

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	11	26	32	17	9	95	0.21	1.14	0.19	0.14
2	8	25	26	24	13	96	0.2	1.15	0.18	0.13
3	11	21	29	14	12	87	0.2	1.16	0.18	0.13
4	9	23	27	21	11	91	0.21	1.21	0.2	0.13
5	11	25	30	16	11	93	0.21	1.12	0.18	0.14
6	6	28	30	26	9	99	0.22	1.14	0.19	0.15
7	8	25	30	17	9	89	0.21	1.15	0.19	0.13
8	11	27	33	20	12	103	0.2	1.15	0.16	0.13
9	11	17	34	19	11	92	0.22	1.19	0.19	0.15
10	9	33	38	16	13	109	0.2	1.14	0.18	0.13

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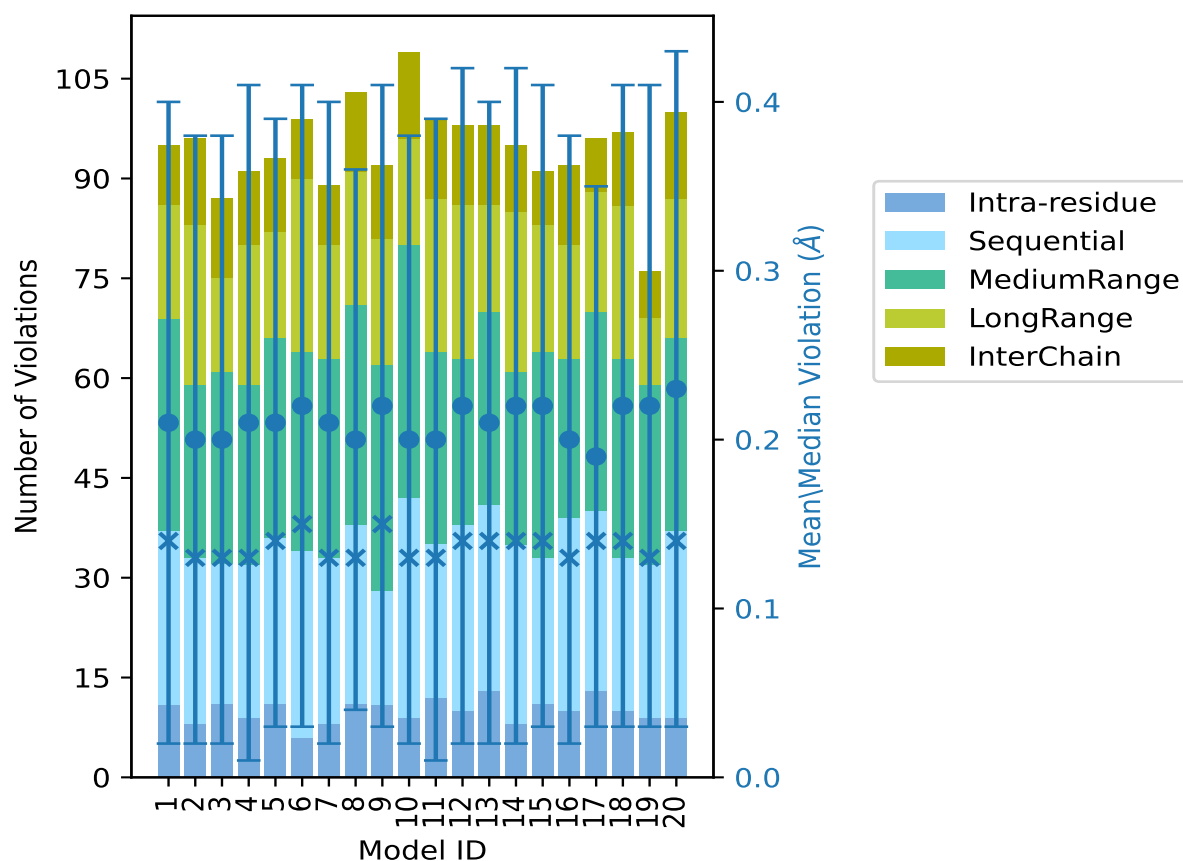
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
11	12	23	29	23	12	99	0.2	1.24	0.19	0.13
12	10	28	25	23	12	98	0.22	1.19	0.2	0.14
13	13	28	29	16	12	98	0.21	1.15	0.19	0.14
14	8	27	26	24	10	95	0.22	1.15	0.2	0.14
15	11	22	31	19	8	91	0.22	1.16	0.19	0.14
16	10	29	24	17	12	92	0.2	1.16	0.18	0.13
17	13	27	30	18	8	96	0.19	1.14	0.16	0.14
18	10	23	30	23	11	97	0.22	1.15	0.19	0.14
19	9	23	27	10	7	76	0.22	1.17	0.19	0.13
20	9	28	29	21	13	100	0.23	1.21	0.2	0.14

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 ⓘ

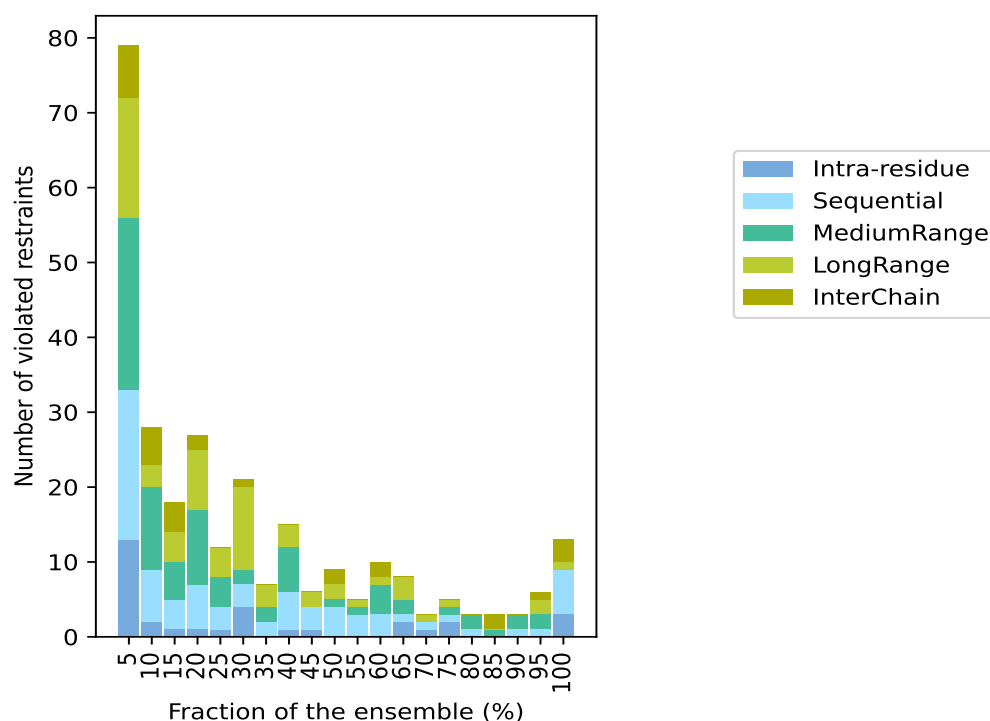
Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 3749(IR:930, SQ:768, MR:1003, LR:897, IC:151) restraints 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	%
13	20	23	16	7	79	1	5.0
2	7	11	3	5	28	2	10.0
1	4	5	4	4	18	3	15.0
1	6	10	8	2	27	4	20.0
1	3	4	4	0	12	5	25.0
4	3	2	11	1	21	6	30.0
0	2	2	3	0	7	7	35.0
1	5	6	3	0	15	8	40.0
1	3	0	2	0	6	9	45.0
0	4	1	2	2	9	10	50.0
0	3	1	1	0	5	11	55.0
0	3	4	1	2	10	12	60.0
2	1	2	3	0	8	13	65.0
1	1	0	1	0	3	14	70.0
2	1	1	1	0	5	15	75.0
0	1	2	0	0	3	16	80.0
0	0	1	0	2	3	17	85.0
0	1	2	0	0	3	18	90.0
0	1	2	2	1	6	19	95.0
3	6	0	1	3	13	20	100.0

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

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

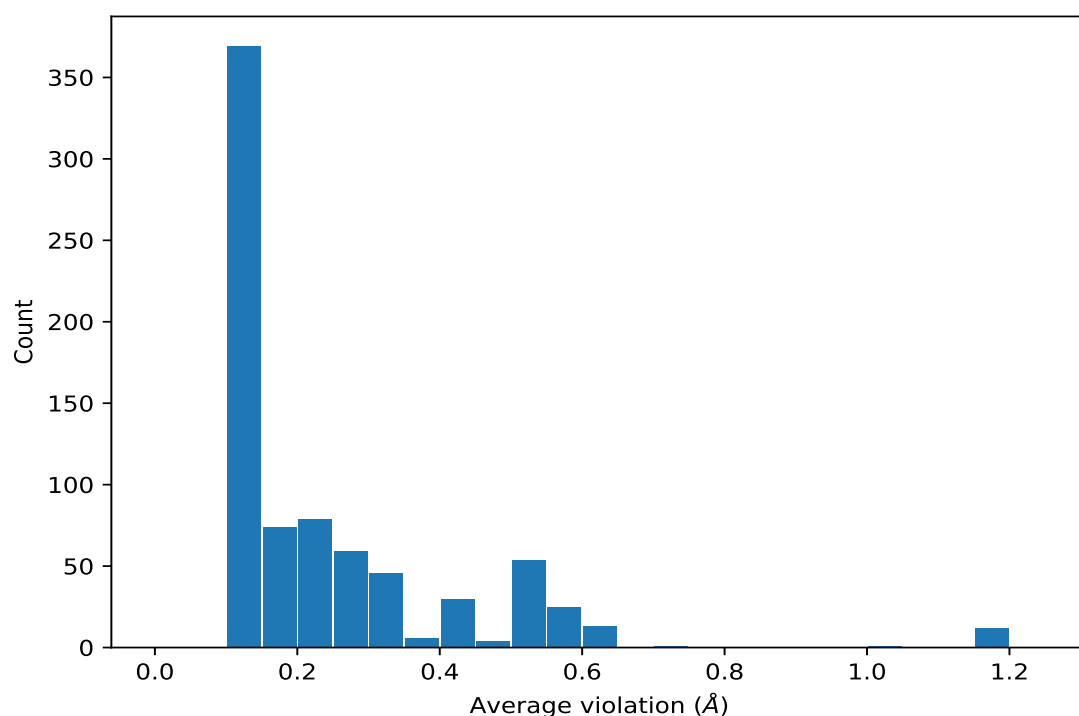
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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 violations for the 10 worst performing restraints, sorted by number of violated models and the mean violation value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	20	1.16	0.03	1.15
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	20	1.16	0.03	1.15
(1,3892)	4:402:B:ATP:PG	4:402:B:ATP:O3A	20	0.74	0.05	0.74
(1,3627)	1:311:B:ARG:HE	1:310:B:SER:HA	20	0.61	0.14	0.57
(1,3584)	1:367:B:TYR:H	1:368:B:LEU:HD11	20	0.57	0.03	0.57
(1,3584)	1:367:B:TYR:H	1:368:B:LEU:HD12	20	0.57	0.03	0.57

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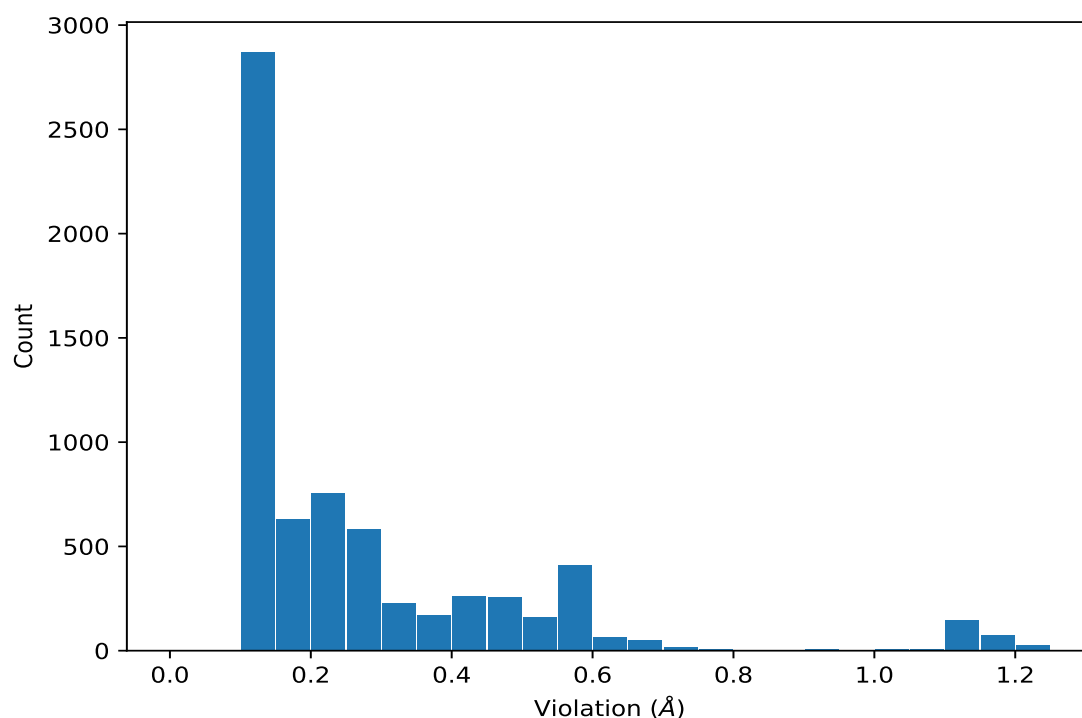
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,3584)	1:367:B:TYR:H	1:368:B:LEU:HD13	20	0.57	0.03	0.57
(1,3584)	1:367:B:TYR:H	1:368:B:LEU:HD21	20	0.57	0.03	0.57
(1,3584)	1:367:B:TYR:H	1:368:B:LEU:HD22	20	0.57	0.03	0.57
(1,3584)	1:367:B:TYR:H	1:368:B:LEU:HD23	20	0.57	0.03	0.57
(1,151)	1:356:B:THR:HG1	2:5:A:DG:H22	20	0.55	0.09	0.53
(1,66)	1:318:B:LEU:HD21	2:6:A:DC:H1'	20	0.45	0.08	0.47
(1,66)	1:318:B:LEU:HD23	2:6:A:DC:H6	20	0.45	0.08	0.47
(1,66)	1:318:B:LEU:HD23	2:6:A:DC:H1'	20	0.45	0.08	0.47
(1,66)	1:318:B:LEU:HD22	2:6:A:DC:H6	20	0.45	0.08	0.47
(1,2939)	1:268:B:VAL:HG11	1:269:B:LYS:HB3	20	0.38	0.05	0.4
(1,2939)	1:268:B:VAL:HG12	1:269:B:LYS:HB3	20	0.38	0.05	0.4
(1,2939)	1:268:B:VAL:HG13	1:269:B:LYS:HB3	20	0.38	0.05	0.4
(1,2939)	1:268:B:VAL:HG21	1:269:B:LYS:HB3	20	0.38	0.05	0.4
(1,2939)	1:268:B:VAL:HG22	1:269:B:LYS:HB3	20	0.38	0.05	0.4
(1,2939)	1:268:B:VAL:HG23	1:269:B:LYS:HB3	20	0.38	0.05	0.4
(1,125)	1:355:B:ILE:HD12	2:2:A:DT:H3'	20	0.31	0.05	0.31
(3,16)	1:266:B:GLU:H	1:262:B:GLU:O	20	0.27	0.02	0.27
(1,3348)	1:332:B:LEU:HD11	1:353:B:PHE:HA	20	0.25	0.09	0.23

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 provides the 10 worst performing restraints, sorted by the violation value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	11	1.24
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	11	1.24
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	11	1.24
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	11	1.24
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	11	1.24
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	11	1.24
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	11	1.24
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	11	1.24
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	11	1.24
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	11	1.24
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	11	1.24
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	11	1.24
(1,3899)	4:402:B:ATP:PG	1:309:B:ARG:NH1	20	1.21
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	4	1.21
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	4	1.21
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	4	1.21
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	4	1.21

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	4	1.21
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	4	1.21
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	4	1.21
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	4	1.21
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	4	1.21
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	4	1.21
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	4	1.21
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	4	1.21
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	9	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	9	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	9	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	9	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	9	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	9	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	9	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	9	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	9	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	9	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	9	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	9	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	12	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	12	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	12	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	12	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	12	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	12	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	12	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	12	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	12	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	12	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	12	1.19
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	12	1.19
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	19	1.17
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	19	1.17
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	19	1.17
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	19	1.17
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	19	1.17
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	19	1.17
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	19	1.17
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	19	1.17
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	19	1.17
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	19	1.17

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	19	1.17
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	19	1.17
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	3	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	3	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	3	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	3	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	3	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	3	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	3	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	3	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	3	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	3	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	3	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	3	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	15	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	15	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	15	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	15	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	15	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	15	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	15	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	15	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	15	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	15	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	15	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	15	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	16	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD12	16	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD13	16	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD21	16	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD22	16	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD23	16	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD11	16	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD12	16	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD13	16	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD21	16	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD22	16	1.16
(1,3587)	1:367:B:TYR:HD2	1:368:B:LEU:HD23	16	1.16
(1,3587)	1:367:B:TYR:HD1	1:368:B:LEU:HD11	2	1.15

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

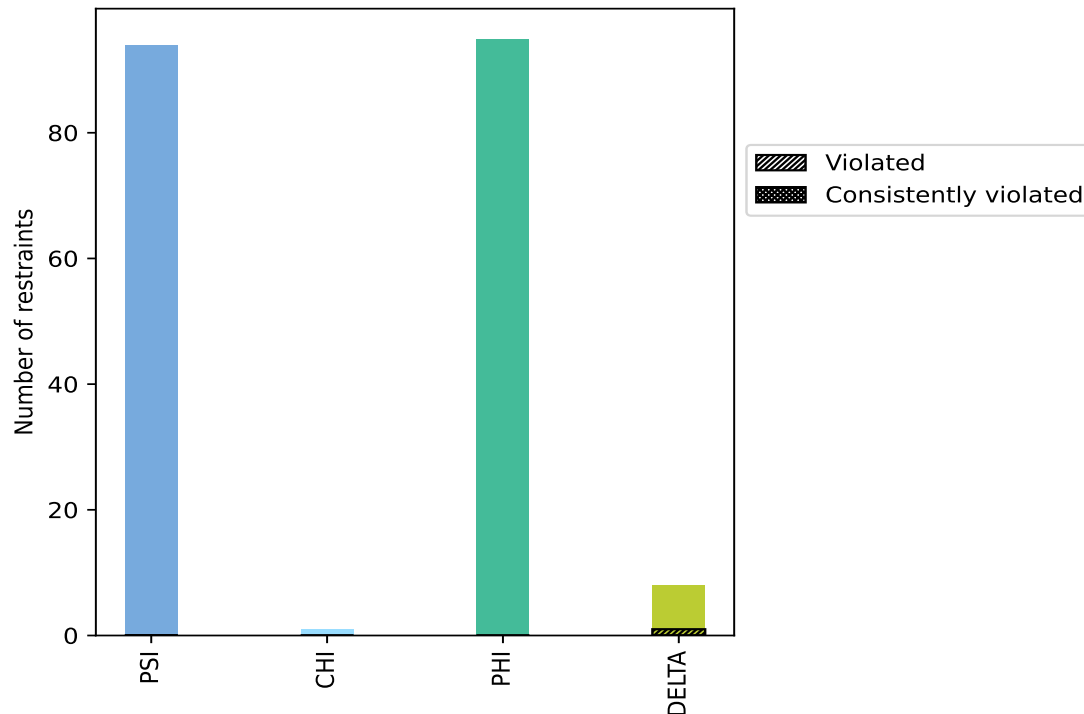
10.1 Summary of dihedral-angle violations [i](#)

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	94	47.5	0	0.0	0.0	0	0.0	0.0
CHI	1	0.5	0	0.0	0.0	0	0.0	0.0
PHI	95	48.0	0	0.0	0.0	0	0.0	0.0
DELTA	8	4.0	1	12.5	0.5	0	0.0	0.0
Total	198	100.0	1	0.5	0.5	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](#)



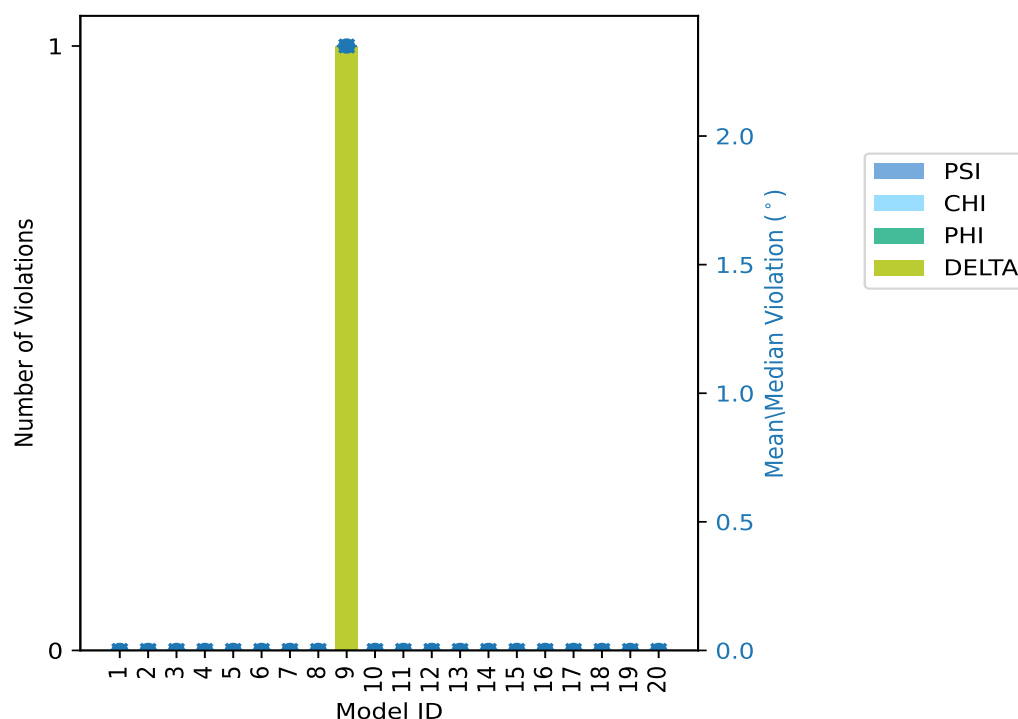
Violated and consistently violated restraints are shown using different hatch patterns in their respective categories

10.2 Dihedral-angle violation statistics for each model

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	CHI	PHI	DELTA	Total				
1	0	0	0	0	0	0.0	0.0	0.0	0.0
2	0	0	0	0	0	0.0	0.0	0.0	0.0
3	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
5	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
7	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
9	0	0	0	1	1	2.35	2.35	0.0	2.35
10	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
12	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
14	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
16	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
18	0	0	0	0	0	0.0	0.0	0.0	0.0
19	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

10.2.1 Bar graph : Dihedral violation statistics for each model [i](#)



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

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

Violation analysis may find that some restraints are violated in very few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of ensemble.

Number of violated restraints					Fraction of the ensemble	
PSI	CHI	PHI	DELTA	Total	Count ¹	%
0	0	0	1	1	1	5.0
0	0	0	0	0	2	10.0
0	0	0	0	0	3	15.0
0	0	0	0	0	4	20.0
0	0	0	0	0	5	25.0
0	0	0	0	0	6	30.0
0	0	0	0	0	7	35.0
0	0	0	0	0	8	40.0
0	0	0	0	0	9	45.0
0	0	0	0	0	10	50.0
0	0	0	0	0	11	55.0

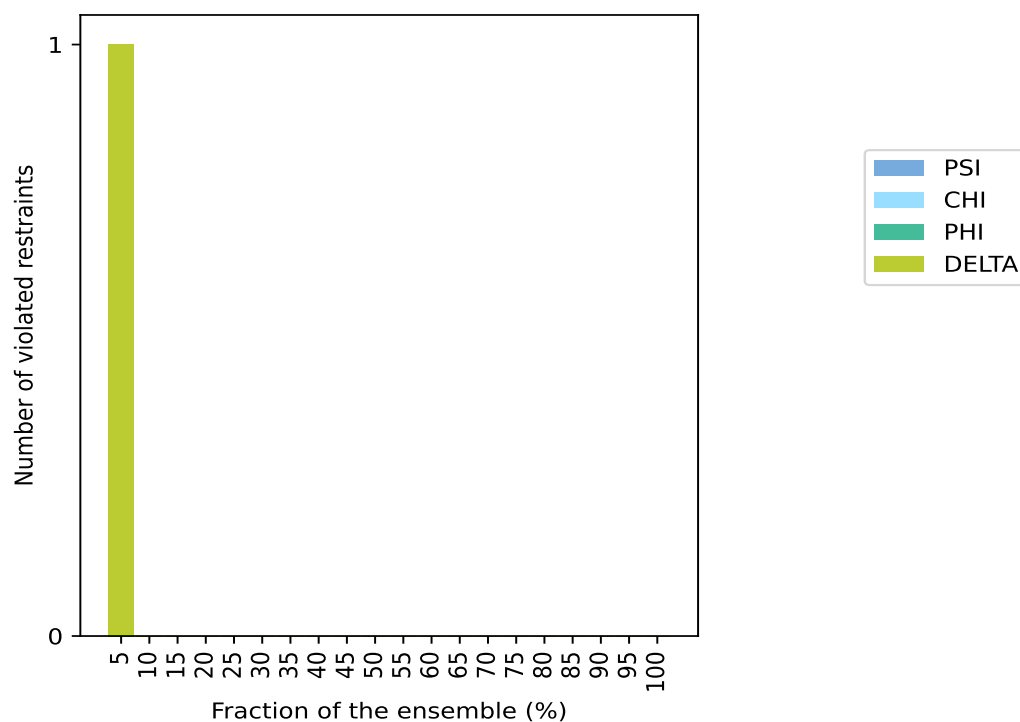
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Number of violated restraints					Fraction of the ensemble	
PSI	CHI	PHI	DELTA	Total	Count ¹	%
0	0	0	0	0	12	60.0
0	0	0	0	0	13	65.0
0	0	0	0	0	14	70.0
0	0	0	0	0	15	75.0
0	0	0	0	0	16	80.0
0	0	0	0	0	17	85.0
0	0	0	0	0	18	90.0
0	0	0	0	0	19	95.0
0	0	0	0	0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

¹ Number of models with violations

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



10.4 Most violated dihedral-angle restraints in the ensemble ⓘ

No violations found

10.5 All violated dihedral-angle restraints [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.

Data insufficient to plot histogram

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 (°)
(1,4)	2:5:A:DG:C5'	2:5:A:DG:C4'	2:5:A:DG:C3'	2:5:A:DG:O3'	9	2.35