



Full wwPDB NMR Structure Validation Report ⓘ

Jun 5, 2023 – 03:36 AM EDT

PDB ID : 2LX1
BMRB ID : 18656
Title : Major Conformation of the Internal Loop 5'GAGU/3'UGAG
Authors : Kennedy, S.D.; Turner, D.H.
Deposited on : 2012-08-12

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 : 20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
BMRB Restraints Analysis : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.33

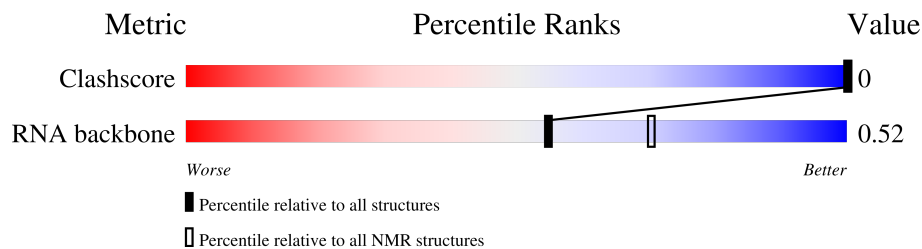
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 28%.

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)
Clashescore	158937	12864
RNA backbone	4643	676

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	11	36% 45% 18%
1	B	11	27% 64% 9%

2 Ensemble composition and analysis

This entry contains 19 models. This entry does not contain polypeptide chains, therefore identification of well-defined residues and clustering analysis are not possible. All residues are included in the validation scores.

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 712 atoms, of which 242 are hydrogens and 0 are deuteriums.

- Molecule 1 is a RNA chain called RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3').

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		P
1	A	11	Total 356	106	121	45	74	10	0
1	B	11	Total 356	106	121	45	74	10	0

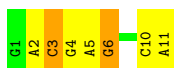
4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

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

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A: 



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain B: 



4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A: 



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain B: 



4.2.2 Score per residue for model 2

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

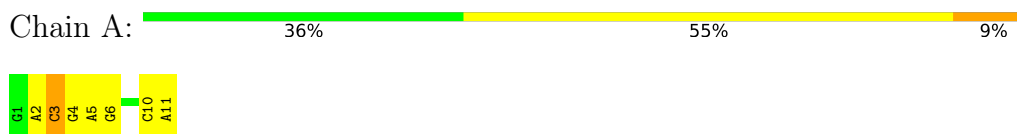


- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')



4.2.3 Score per residue for model 3

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')



4.2.4 Score per residue for model 4

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')



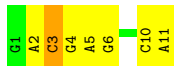
- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')



4.2.5 Score per residue for model 5

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  36% 55% 9%



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain B:  27% 45% 27%




4.2.6 Score per residue for model 6

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  36% 55% 9%



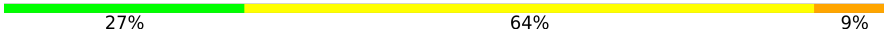
- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

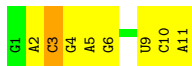
Chain B:  36% 55% 9%



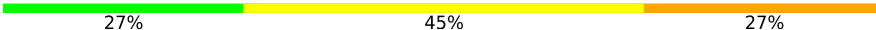
4.2.7 Score per residue for model 7

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  27% 64% 9%



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

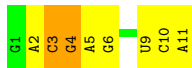
Chain B:  27% 45% 27%



4.2.8 Score per residue for model 8

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A: 27% 55% 18%



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

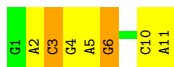
Chain B: 36% 45% 18%



4.2.9 Score per residue for model 9

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A: 36% 45% 18%



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain B: 18% 64% 18%



4.2.10 Score per residue for model 10

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A: 27% 64% 9%



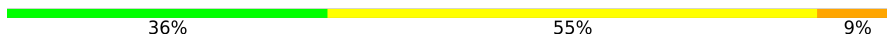
- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

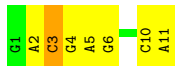
Chain B: 27% 64% 9%



4.2.11 Score per residue for model 11

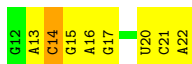
- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  36% 55% 9%



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

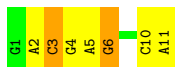
Chain B:  27% 64% 9%



4.2.12 Score per residue for model 12

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  36% 45% 18%



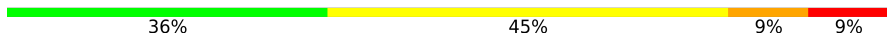
- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain B:  27% 64% 9%



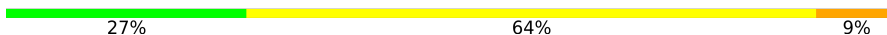
4.2.13 Score per residue for model 13

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  36% 45% 9% 9%



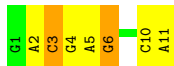
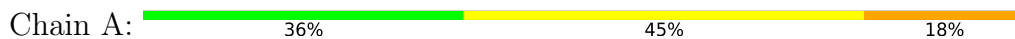
- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain B:  27% 64% 9%



4.2.14 Score per residue for model 14

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

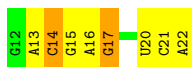
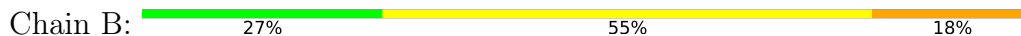


4.2.15 Score per residue for model 15

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

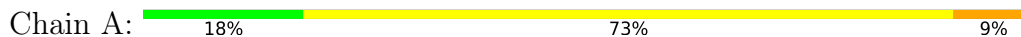


- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

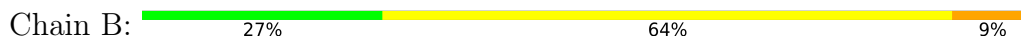


4.2.16 Score per residue for model 16

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

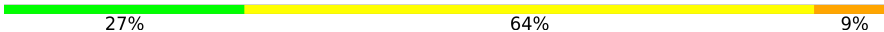


- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')



4.2.17 Score per residue for model 17

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  27% 64% 9%



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

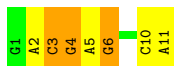
Chain B:  36% 55% 9%




4.2.18 Score per residue for model 18

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  36% 36% 27%



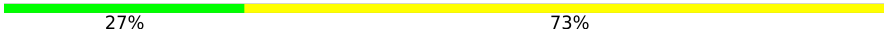
- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

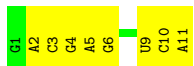
Chain B:  36% 55% 9%



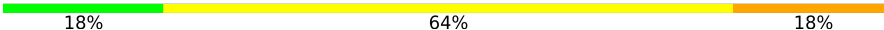
4.2.19 Score per residue for model 19

- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain A:  27% 73%



- Molecule 1: RNA (5'-R(*GP*AP*CP*GP*AP*GP*UP*GP*UP*CP*A)-3')

Chain B:  18% 64% 18%



5 Refinement protocol and experimental data overview

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

Of the 600 calculated structures, 19 were deposited, based on the following criterion: *structures with the least restraint violations*.

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

Software name	Classification	Version
Amber	structure solution	10
Amber	refinement	10

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	220
Number of shifts mapped to atoms	220
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	28%

6 Model quality i

6.1 Standard geometry i

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	A	1.40±0.01	0±0/263 (0.0± 0.0%)	2.21±0.01	16±1/409 (4.0± 0.2%)
1	B	1.40±0.01	0±0/263 (0.0± 0.0%)	2.22±0.02	16±1/409 (4.0± 0.2%)
All	All	1.40	0/9994 (0.0%)	2.22	624/15542 (4.0%)

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

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	1.2±0.6
1	B	0.0±0.0	1.6±0.6
All	All	0	52

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	16	A	N1-C6-N6	-9.96	112.62	118.60	4	19
1	A	5	A	N1-C6-N6	-9.92	112.65	118.60	9	19
1	A	11	A	N1-C6-N6	-9.57	112.86	118.60	6	19
1	B	22	A	N1-C6-N6	-9.27	113.04	118.60	3	19
1	A	2	A	N1-C6-N6	-8.85	113.29	118.60	2	19
1	B	13	A	N1-C6-N6	-8.80	113.32	118.60	5	19
1	B	22	A	C5-C6-N1	8.10	121.75	117.70	3	19
1	A	11	A	C5-C6-N1	8.07	121.74	117.70	13	19
1	B	16	A	C5-C6-N1	8.01	121.70	117.70	5	19
1	A	5	A	C5-C6-N1	7.84	121.62	117.70	7	19
1	B	14	C	N3-C2-O2	-7.49	116.66	121.90	4	19
1	A	10	C	N3-C2-O2	-7.34	116.76	121.90	14	19
1	B	21	C	N3-C2-O2	-7.30	116.79	121.90	1	19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	B	15	G	O4'-C1'-N9	7.25	114.00	108.20	1	19
1	A	2	A	C5-C6-N1	7.24	121.32	117.70	4	19
1	A	3	C	N3-C2-O2	-7.21	116.86	121.90	10	19
1	A	4	G	O4'-C1'-N9	7.12	113.90	108.20	8	19
1	A	11	A	C4-C5-C6	-7.08	113.46	117.00	13	19
1	B	13	A	C5-C6-N1	7.06	121.23	117.70	4	19
1	B	22	A	C4-C5-C6	-6.92	113.54	117.00	13	19
1	A	5	A	C4-C5-C6	-6.86	113.57	117.00	4	19
1	B	16	A	C4-C5-C6	-6.79	113.61	117.00	4	19
1	A	2	A	C4-C5-C6	-6.50	113.75	117.00	2	19
1	B	13	A	C4-C5-C6	-6.46	113.77	117.00	2	19
1	A	10	C	O4'-C1'-N1	6.11	113.09	108.20	4	16
1	B	21	C	N1-C2-O2	6.07	122.54	118.90	1	19
1	B	15	G	N3-C4-C5	-6.00	125.60	128.60	15	16
1	B	14	C	N1-C2-O2	6.00	122.50	118.90	7	19
1	A	10	C	N1-C2-O2	5.94	122.47	118.90	7	18
1	B	21	C	O4'-C1'-N1	5.94	112.95	108.20	2	16
1	A	4	G	N3-C4-C5	-5.83	125.69	128.60	7	14
1	A	3	C	N1-C2-O2	5.79	122.37	118.90	18	19
1	A	6	G	N3-C4-C5	-5.62	125.79	128.60	14	10
1	B	17	G	N3-C4-C5	-5.54	125.83	128.60	8	9
1	A	9	U	O4'-C1'-N1	5.53	112.62	108.20	16	5
1	B	20	U	O4'-C1'-N1	5.41	112.53	108.20	11	3
1	B	19	G	N1-C6-O6	-5.38	116.67	119.90	9	3
1	A	6	G	N1-C6-O6	-5.02	116.89	119.90	14	1

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	B	14	C	Sidechain	18
1	A	3	C	Sidechain	15
1	B	19	G	Sidechain	9
1	A	4	G	Sidechain	3
1	B	15	G	Sidechain	3
1	A	8	G	Sidechain	2
1	A	2	A	Sidechain	1
1	A	5	A	Sidechain	1

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
All	All	8930	4598	4598	-

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 -.

There are no clashes.

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

There are no protein molecules in this entry.

6.3.2 Protein sidechains [i](#)

There are no protein molecules in this entry.

6.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers	Suiteness
1	A	10/11 (91%)	1±0 (11±2%)	0±0 (1±3%)	0.51±0.04
1	B	10/11 (91%)	1±0 (11±2%)	0±0 (0±0%)	0.54±0.04
All	All	380/418 (91%)	40 (11%)	2 (1%)	0.52

The overall RNA backbone suiteness is 0.52.

All unique RNA backbone outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
1	A	6	G	19
1	B	17	G	19
1	B	18	U	1
1	A	7	U	1

All unique RNA pucker outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
1	A	4	G	2

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation i

The completeness of assignment taking into account all chemical shift lists is 28% for the well-defined parts and 28% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_3*

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	106
Number of shifts mapped to atoms	106
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	2

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 22%, i.e. 94 atoms were assigned a chemical shift out of a possible 418. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Sugar	66/242 (27%)	66/132 (50%)	0/110 (0%)	0/0 (—%)
Base	28/176 (16%)	15/110 (14%)	13/36 (36%)	0/30 (0%)
Overall	94/418 (22%)	81/242 (33%)	13/146 (9%)	0/30 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 22%, i.e. 94 atoms were assigned a chemical shift out of a possible 418. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Sugar	66/242 (27%)	66/132 (50%)	0/110 (0%)	0/0 (—%)
Base	28/176 (16%)	15/110 (14%)	13/36 (36%)	0/30 (0%)
Overall	94/418 (22%)	81/242 (33%)	13/146 (9%)	0/30 (0%)

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	5	A	H5''	2.85	3.15 – 5.18	-6.5
1	A	6	G	H2'	6.14	3.35 – 5.81	6.3

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_chem_shift_list_3_dup*

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	21
Number of shifts mapped to atoms	21
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

7.2.2 Chemical shift referencing [i](#)

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

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 4%, i.e. 15 atoms were assigned a chemical shift out of a possible 418. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Sugar	0/242 (0%)	0/132 (0%)	0/110 (0%)	0/0 (—%)
Base	15/176 (9%)	15/110 (14%)	0/36 (0%)	0/30 (0%)
Overall	15/418 (4%)	15/242 (6%)	0/146 (0%)	0/30 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 4%, i.e. 15 atoms were assigned a chemical shift out of a possible 418. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Sugar	0/242 (0%)	0/132 (0%)	0/110 (0%)	0/0 (—%)
Base	15/176 (9%)	15/110 (14%)	0/36 (0%)	0/30 (0%)
Overall	15/418 (4%)	15/242 (6%)	0/146 (0%)	0/30 (0%)

7.2.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.2.5 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

7.3 Chemical shift list 3

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

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	47
Number of shifts mapped to atoms	47

Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

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 11%, i.e. 45 atoms were assigned a chemical shift out of a possible 418. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Sugar	29/242 (12%)	29/132 (22%)	0/110 (0%)	0/0 (—%)
Base	16/176 (9%)	16/110 (15%)	0/36 (0%)	0/30 (0%)
Overall	45/418 (11%)	45/242 (19%)	0/146 (0%)	0/30 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 11%, i.e. 45 atoms were assigned a chemical shift out of a possible 418. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Sugar	29/242 (12%)	29/132 (22%)	0/110 (0%)	0/0 (—%)
Base	16/176 (9%)	16/110 (15%)	0/36 (0%)	0/30 (0%)
Overall	45/418 (11%)	45/242 (19%)	0/146 (0%)	0/30 (0%)

7.3.4 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
3	A	5	A	H5''	2.80	3.15 – 5.18	-6.7

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: *assigned_chem_shift_list_2*

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	46
Number of shifts mapped to atoms	46
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 11%, i.e. 44 atoms were assigned a chemical shift out of a possible 418. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Sugar	10/242 (4%)	10/132 (8%)	0/110 (0%)	0/0 (—%)
Base	34/176 (19%)	29/110 (26%)	0/36 (0%)	5/30 (17%)
Overall	44/418 (11%)	39/242 (16%)	0/146 (0%)	5/30 (17%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 11%, i.e. 44 atoms were assigned a chemical shift out of a possible 418. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Sugar	10/242 (4%)	10/132 (8%)	0/110 (0%)	0/0 (—%)
Base	34/176 (19%)	29/110 (26%)	0/36 (0%)	5/30 (17%)
Overall	44/418 (11%)	39/242 (16%)	0/146 (0%)	5/30 (17%)

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	167
Intra-residue ($ i-j =0$)	74
Sequential ($ i-j =1$)	52
Medium range ($ i-j >1$ and $ i-j <5$)	10
Long range ($ i-j \geq 5$)	31
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	7.6
Number of long range restraints per residue ¹	1.4

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)	4.1	0.2
0.2-0.5 (Medium)	1.6	0.28
>0.5 (Large)	45.6	33.39

8.2.2 Average number of dihedral-angle violations per model

Dihedral-angle violations less than 1° are not included in the calculation. There are no dihedral-angle violations

9 Distance violation analysis [i](#)

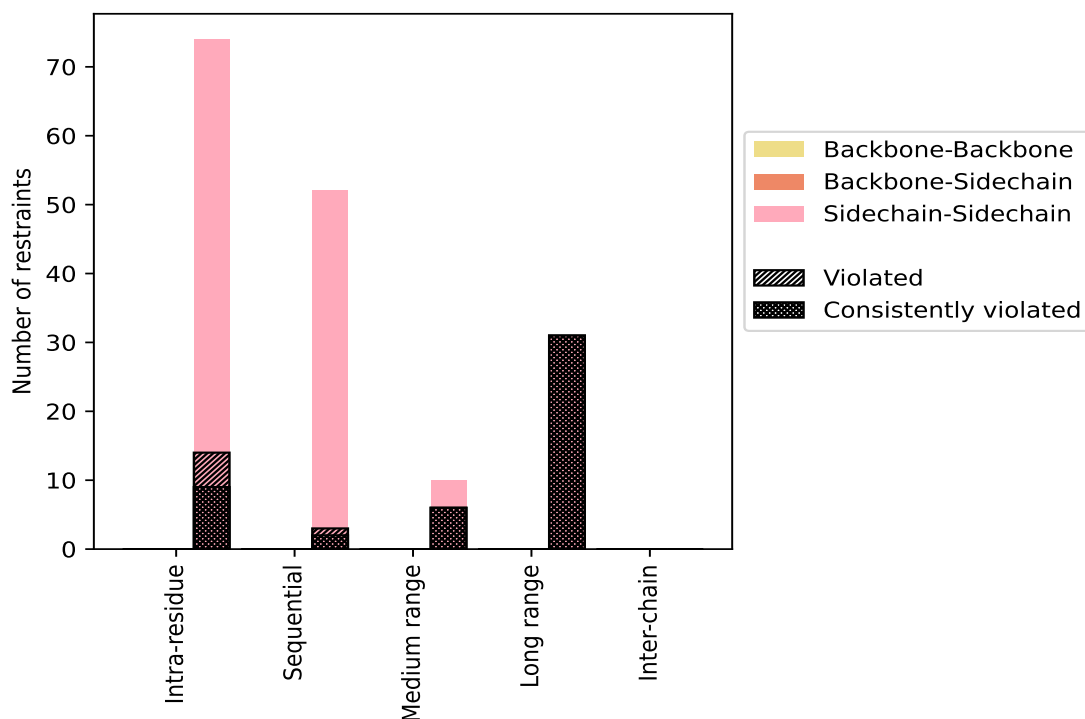
9.1 Summary of distance violations [i](#)

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

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	74	44.3	14	18.9	8.4	9	12.2	5.4
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	74	44.3	14	18.9	8.4	9	12.2	5.4
Sequential ($i-j =1$)	52	31.1	3	5.8	1.8	2	3.8	1.2
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	52	31.1	3	5.8	1.8	2	3.8	1.2
Medium range ($i-j >1$ & $i-j <5$)	10	6.0	6	60.0	3.6	6	60.0	3.6
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	10	6.0	6	60.0	3.6	6	60.0	3.6
Long range ($i-j \geq 5$)	31	18.6	31	100.0	18.6	31	100.0	18.6
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	31	18.6	31	100.0	18.6	31	100.0	18.6
Inter-chain	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Hydrogen bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	167	100.0	54	32.3	32.3	48	28.7	28.7
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	167	100.0	54	32.3	32.3	48	28.7	28.7

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 disulfid bonds are counted in their appropriate category on the x-axis

9.2 Distance violation statistics for each model [i](#)

The following table provides the distance violation statistics for each model in the ensemble. Violations less than 0.1 Å are not included in the statistics.

Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	12	3	6	31	0	52	14.72	31.78	11.24	15.67
2	12	3	6	31	0	52	14.76	31.67	11.27	15.76
3	12	3	6	31	0	52	14.75	31.39	11.32	15.75
4	10	3	6	31	0	50	15.25	32.25	11.22	16.04
5	14	2	6	31	0	53	14.53	32.5	11.41	14.87
6	12	2	6	31	0	51	15.15	32.56	11.34	16.75
7	12	3	6	31	0	52	14.82	32.57	11.45	15.53
8	11	2	6	31	0	50	15.36	32.15	11.24	16.37
9	12	3	6	31	0	52	14.94	33.39	11.56	16.01
10	13	2	6	31	0	52	14.89	31.81	11.34	15.89
11	12	2	6	31	0	51	15.11	32.11	11.32	16.76

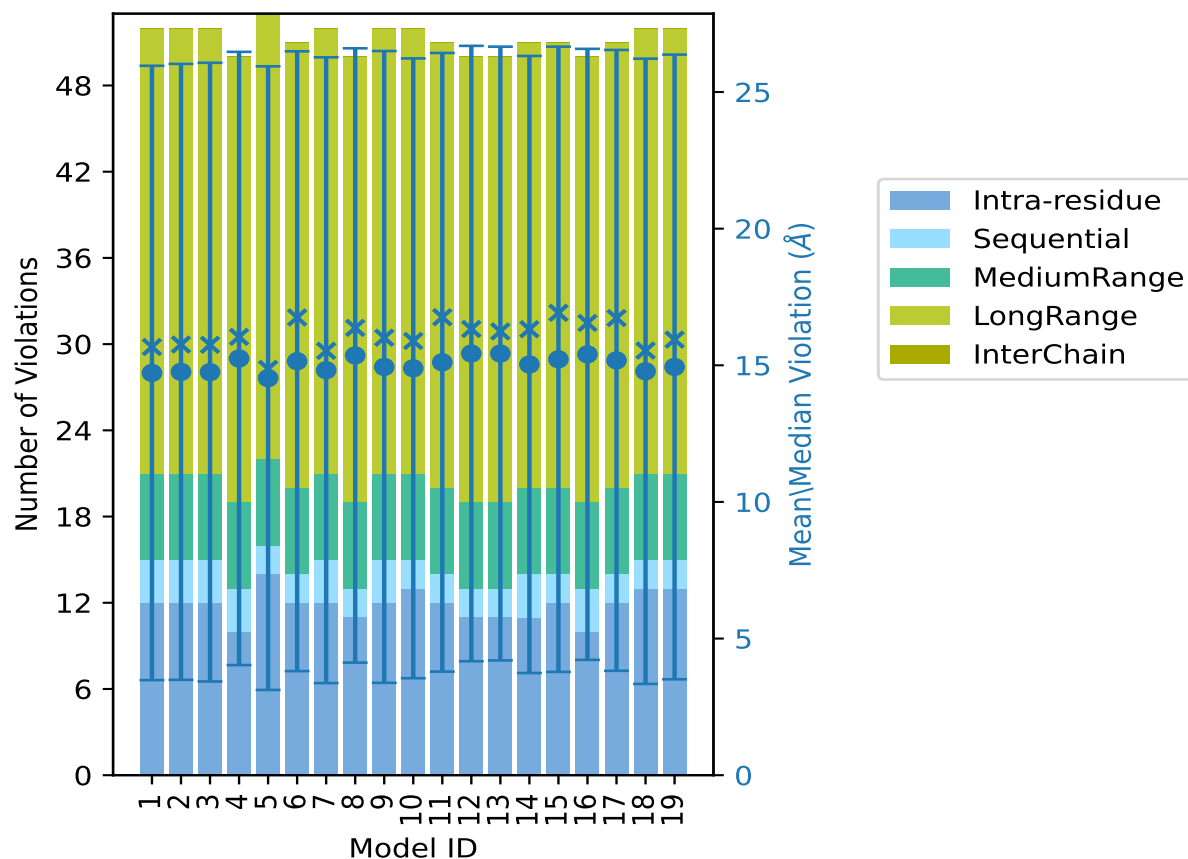
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
12	11	2	6	31	0	50	15.43	32.03	11.26	16.33
13	11	2	6	31	0	50	15.43	31.59	11.23	16.24
14	11	3	6	31	0	51	15.03	32.17	11.29	16.32
15	12	2	6	31	0	51	15.22	32.77	11.44	16.92
16	10	3	6	31	0	50	15.4	32.1	11.18	16.55
17	12	2	6	31	0	51	15.18	32.42	11.36	16.73
18	13	2	6	31	0	52	14.78	32.48	11.44	15.54
19	13	2	6	31	0	52	14.94	32.65	11.43	15.94

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 [i](#)



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

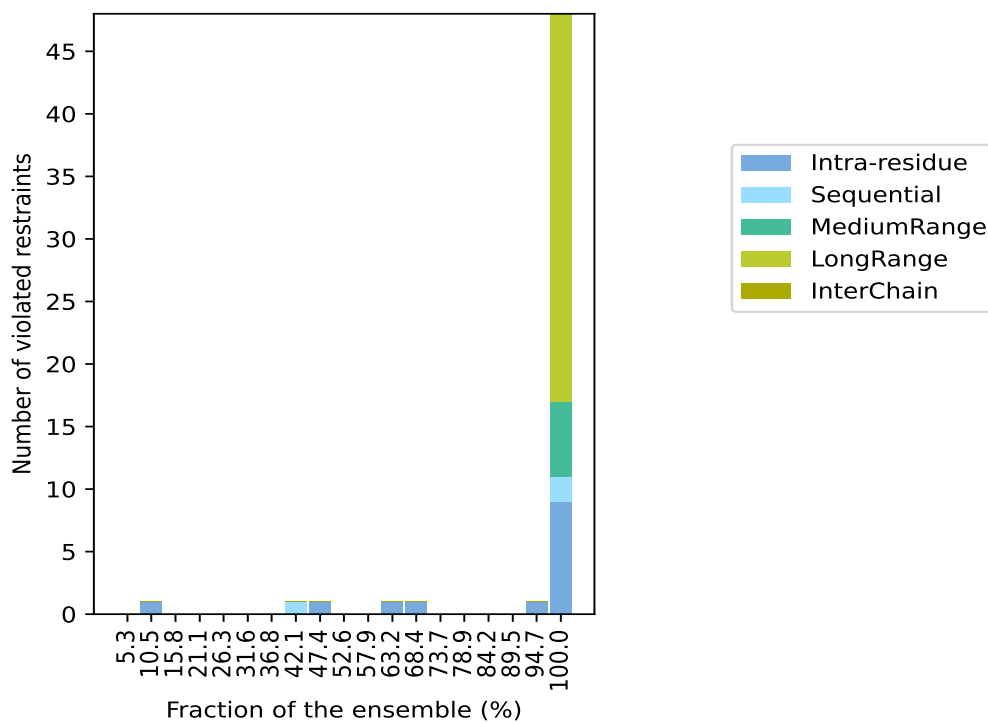
9.3 Distance violation statistics for the ensemble

Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 113(IR:60, SQ:49, MR:4, LR:0, 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	0	0	0	0	0	1	5.3
1	0	0	0	0	1	2	10.5
0	0	0	0	0	0	3	15.8
0	0	0	0	0	0	4	21.1
0	0	0	0	0	0	5	26.3
0	0	0	0	0	0	6	31.6
0	0	0	0	0	0	7	36.8
0	1	0	0	0	1	8	42.1
1	0	0	0	0	1	9	47.4
0	0	0	0	0	0	10	52.6
0	0	0	0	0	0	11	57.9
1	0	0	0	0	1	12	63.2
1	0	0	0	0	1	13	68.4
0	0	0	0	0	0	14	73.7
0	0	0	0	0	0	15	78.9
0	0	0	0	0	0	16	84.2
0	0	0	0	0	0	17	89.5
1	0	0	0	0	1	18	94.7
9	2	6	31	0	48	19	100.0

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

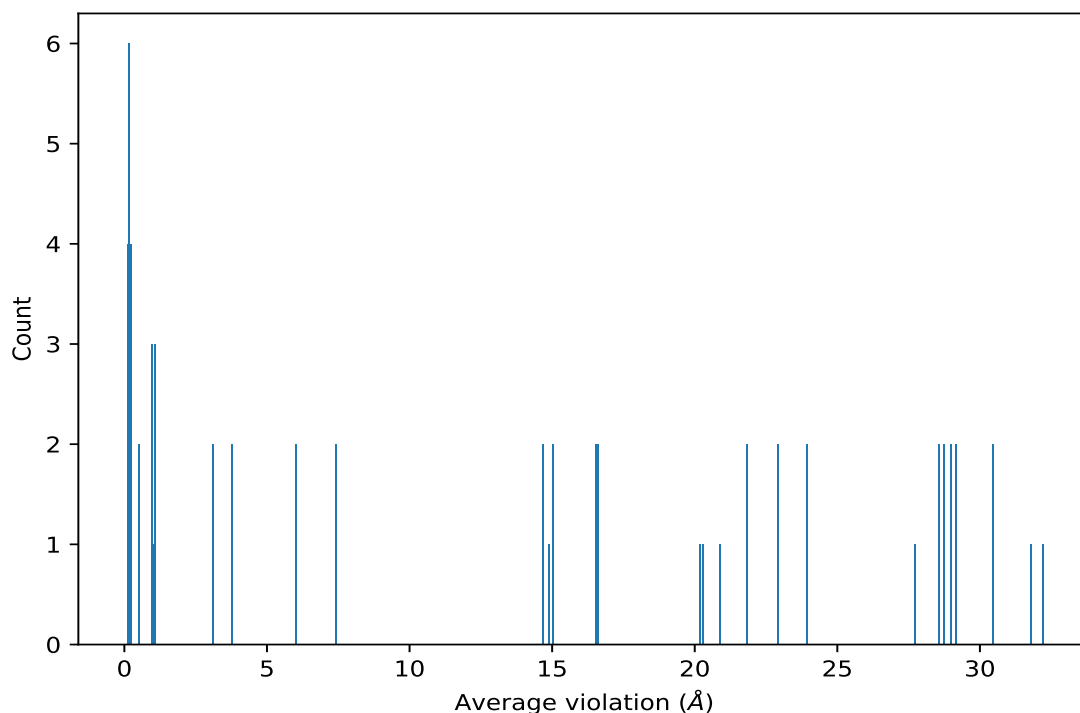
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



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)
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	19	32.23	0.46	32.17
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	19	31.78	0.46	31.72
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	19	30.45	0.34	30.49
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	19	30.45	0.34	30.49
(3,3)	1:B:12:G:H22	1:B:21:C:O2	19	29.18	0.18	29.2
(3,16)	1:B:21:C:O2	1:B:12:G:H22	19	29.18	0.18	29.2
(3,2)	1:B:12:G:O6	1:B:21:C:H42	19	28.96	0.21	28.95
(3,15)	1:B:21:C:H42	1:B:12:G:O6	19	28.96	0.21	28.95
(1,3)	1:B:12:G:H1	1:B:22:A:H2	19	28.72	0.33	28.75
(1,30)	1:B:22:A:H2	1:B:12:G:H1	19	28.72	0.33	28.75
(3,1)	1:B:12:G:H1	1:B:21:C:N3	19	28.56	0.1	28.54
(3,14)	1:B:21:C:N3	1:B:12:G:H1	19	28.56	0.1	28.54
(1,2)	1:B:12:G:H1	1:B:21:C:H42	19	27.7	0.18	27.72
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	19	23.93	0.22	23.94
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	19	23.93	0.22	23.94
(3,5)	1:B:13:A:H62	1:B:20:U:O4	19	22.9	0.24	22.89

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Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(3,13)	1:B:20:U:O4	1:B:13:A:H62	19	22.9	0.24	22.89
(3,4)	1:B:13:A:N1	1:B:20:U:H3	19	21.83	0.13	21.86
(3,12)	1:B:20:U:H3	1:B:13:A:N1	19	21.83	0.13	21.86
(1,23)	1:B:20:U:H3	1:B:12:G:H1	19	20.88	0.13	20.89
(1,24)	1:B:20:U:H3	1:B:13:A:H2	19	20.27	0.14	20.25
(1,1)	1:B:12:G:H1	1:B:20:U:H3	19	20.16	0.13	20.17
(1,9)	1:B:14:C:H42	1:B:20:U:H3	19	16.62	0.23	16.65
(1,25)	1:B:20:U:H3	1:B:14:C:H42	19	16.62	0.23	16.65
(3,7)	1:B:14:C:H42	1:B:19:G:O6	19	16.5	0.24	16.49
(3,10)	1:B:19:G:O6	1:B:14:C:H42	19	16.5	0.24	16.49
(3,6)	1:B:14:C:N3	1:B:19:G:H1	19	15.01	0.13	14.98
(3,9)	1:B:19:G:H1	1:B:14:C:N3	19	15.01	0.13	14.98
(1,19)	1:B:19:G:H1	1:B:14:C:H42	19	14.88	0.25	14.95
(3,8)	1:B:14:C:O2	1:B:19:G:H22	19	14.65	0.18	14.65
(3,11)	1:B:19:G:H22	1:B:14:C:O2	19	14.65	0.18	14.65
(1,13)	1:B:15:G:H8	1:B:19:G:H1	19	7.42	0.18	7.38
(1,20)	1:B:19:G:H1	1:B:15:G:H8	19	7.42	0.18	7.38
(1,32)	1:B:15:G:H8	1:B:17:G:H1	19	6.02	0.16	6.06
(1,34)	1:B:17:G:H1	1:B:15:G:H8	19	6.02	0.16	6.06
(1,31)	1:B:15:G:H1	1:B:16:A:H2	19	3.75	0.29	3.72
(1,33)	1:B:16:A:H2	1:B:15:G:H1	19	3.75	0.29	3.72
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	19	3.14	0.14	3.12
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	19	3.14	0.14	3.12
(1,142)	1:B:22:A:N7	1:B:22:A:H61	19	1.08	0.0	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	19	1.08	0.01	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	19	1.08	0.01	1.08
(1,139)	1:B:21:C:H5	1:B:21:C:H41	19	1.0	0.01	1.0
(1,138)	1:B:14:C:H5	1:B:14:C:H41	19	0.98	0.01	0.98
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	19	0.53	0.0	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	19	0.53	0.0	0.53
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	19	0.23	0.02	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	19	0.23	0.02	0.23
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	19	0.21	0.05	0.22
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	19	0.21	0.05	0.22
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	19	0.17	0.01	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	19	0.17	0.01	0.17
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	18	0.14	0.01	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	18	0.14	0.01	0.14
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	13	0.13	0.02	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	13	0.13	0.02	0.13
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	12	0.15	0.03	0.15
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	12	0.15	0.03	0.15

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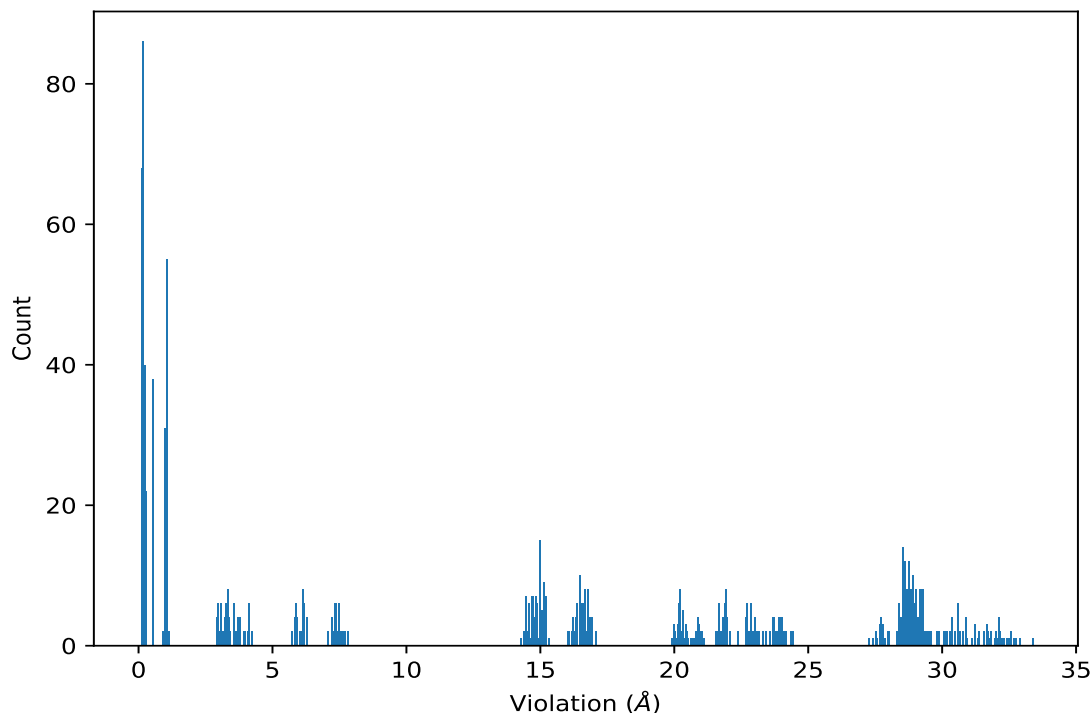
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,135)	1:B:15:G:H1	1:B:15:G:H22	9	0.95	0.0	0.95
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	8	0.15	0.02	0.14
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	8	0.15	0.02	0.14
(1,136)	1:B:17:G:H1	1:B:17:G:H22	2	0.95	0.0	0.95

¹Number of violated models, ²Standard deviation

9.5 All violated distance restraints [i](#)

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

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



9.5.2 Table : All distance violations [i](#)

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

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	9	33.39
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	9	32.94
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	15	32.77
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	19	32.65
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	7	32.57
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	6	32.56
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	5	32.5
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	18	32.48
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	17	32.42
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	15	32.32
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	4	32.25
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	19	32.2
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	14	32.17
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	8	32.15
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	7	32.12
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	11	32.11
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	6	32.11
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	16	32.1
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	5	32.05
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	12	32.03
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	18	32.03
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	17	31.97
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	10	31.81
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	4	31.8
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	1	31.78
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	14	31.72
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	8	31.7
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	2	31.67
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	11	31.66
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	16	31.65
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	13	31.59
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	12	31.58
(1,38)	1:B:12:G:H1'	1:B:22:A:H2	3	31.39
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	10	31.36
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	1	31.33
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	7	31.22
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	7	31.22
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	2	31.22
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	13	31.14
(1,124)	1:B:22:A:H2	1:B:12:G:H1'	3	30.94
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	13	30.86
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	13	30.86
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	9	30.85

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	9	30.85
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	12	30.78
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	12	30.78
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	17	30.69
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	17	30.69
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	18	30.63
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	18	30.63
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	19	30.58
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	19	30.58
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	4	30.55
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	14	30.55
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	4	30.55
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	14	30.55
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	8	30.49
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	8	30.49
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	6	30.35
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	16	30.35
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	6	30.35
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	16	30.35
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	15	30.32
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	15	30.32
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	5	30.29
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	5	30.29
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	11	30.17
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	11	30.17
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	10	30.14
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	10	30.14
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	3	30.07
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	3	30.07
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	2	29.89
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	2	29.89
(1,4)	1:B:12:G:H22	1:B:22:A:H1'	1	29.83
(1,29)	1:B:22:A:H1'	1:B:12:G:H22	1	29.83
(3,3)	1:B:12:G:H22	1:B:21:C:O2	12	29.58
(3,16)	1:B:21:C:O2	1:B:12:G:H22	12	29.58
(1,30)	1:B:22:A:H2	1:B:12:G:H1	9	29.53
(1,3)	1:B:12:G:H1	1:B:22:A:H2	9	29.53
(3,3)	1:B:12:G:H22	1:B:21:C:O2	18	29.45
(3,16)	1:B:21:C:O2	1:B:12:G:H22	18	29.45
(3,2)	1:B:12:G:O6	1:B:21:C:H42	15	29.44
(3,15)	1:B:21:C:H42	1:B:12:G:O6	15	29.44
(3,3)	1:B:12:G:H22	1:B:21:C:O2	13	29.35

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,16)	1:B:21:C:O2	1:B:12:G:H22	13	29.35
(3,3)	1:B:12:G:H22	1:B:21:C:O2	7	29.32
(3,16)	1:B:21:C:O2	1:B:12:G:H22	7	29.32
(3,3)	1:B:12:G:H22	1:B:21:C:O2	14	29.29
(3,16)	1:B:21:C:O2	1:B:12:G:H22	14	29.29
(3,3)	1:B:12:G:H22	1:B:21:C:O2	16	29.27
(3,16)	1:B:21:C:O2	1:B:12:G:H22	16	29.27
(3,2)	1:B:12:G:O6	1:B:21:C:H42	17	29.26
(3,15)	1:B:21:C:H42	1:B:12:G:O6	17	29.26
(3,3)	1:B:12:G:H22	1:B:21:C:O2	10	29.25
(3,16)	1:B:21:C:O2	1:B:12:G:H22	10	29.25
(3,3)	1:B:12:G:H22	1:B:21:C:O2	9	29.22
(3,2)	1:B:12:G:O6	1:B:21:C:H42	11	29.22
(3,16)	1:B:21:C:O2	1:B:12:G:H22	9	29.22
(3,15)	1:B:21:C:H42	1:B:12:G:O6	11	29.22
(3,3)	1:B:12:G:H22	1:B:21:C:O2	4	29.21
(3,16)	1:B:21:C:O2	1:B:12:G:H22	4	29.21
(3,3)	1:B:12:G:H22	1:B:21:C:O2	19	29.2
(3,16)	1:B:21:C:O2	1:B:12:G:H22	19	29.2
(3,3)	1:B:12:G:H22	1:B:21:C:O2	6	29.19
(3,16)	1:B:21:C:O2	1:B:12:G:H22	6	29.19
(3,3)	1:B:12:G:H22	1:B:21:C:O2	15	29.17
(3,2)	1:B:12:G:O6	1:B:21:C:H42	9	29.17
(3,16)	1:B:21:C:O2	1:B:12:G:H22	15	29.17
(3,15)	1:B:21:C:H42	1:B:12:G:O6	9	29.17
(3,2)	1:B:12:G:O6	1:B:21:C:H42	19	29.16
(3,15)	1:B:21:C:H42	1:B:12:G:O6	19	29.16
(3,3)	1:B:12:G:H22	1:B:21:C:O2	5	29.14
(3,16)	1:B:21:C:O2	1:B:12:G:H22	5	29.14
(3,2)	1:B:12:G:O6	1:B:21:C:H42	10	29.1
(3,15)	1:B:21:C:H42	1:B:12:G:O6	10	29.1
(3,3)	1:B:12:G:H22	1:B:21:C:O2	8	29.09
(3,16)	1:B:21:C:O2	1:B:12:G:H22	8	29.09
(1,30)	1:B:22:A:H2	1:B:12:G:H1	7	29.09
(1,3)	1:B:12:G:H1	1:B:22:A:H2	7	29.09
(3,2)	1:B:12:G:O6	1:B:21:C:H42	2	29.03
(3,15)	1:B:21:C:H42	1:B:12:G:O6	2	29.03
(1,30)	1:B:22:A:H2	1:B:12:G:H1	6	29.03
(1,3)	1:B:12:G:H1	1:B:22:A:H2	6	29.03
(3,3)	1:B:12:G:H22	1:B:21:C:O2	3	29.0
(3,3)	1:B:12:G:H22	1:B:21:C:O2	17	29.0
(3,16)	1:B:21:C:O2	1:B:12:G:H22	3	29.0

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,16)	1:B:21:C:O2	1:B:12:G:H22	17	29.0
(3,2)	1:B:12:G:O6	1:B:21:C:H42	6	28.98
(3,2)	1:B:12:G:O6	1:B:21:C:H42	8	28.98
(3,15)	1:B:21:C:H42	1:B:12:G:O6	6	28.98
(3,15)	1:B:21:C:H42	1:B:12:G:O6	8	28.98
(3,2)	1:B:12:G:O6	1:B:21:C:H42	3	28.95
(3,15)	1:B:21:C:H42	1:B:12:G:O6	3	28.95
(3,3)	1:B:12:G:H22	1:B:21:C:O2	11	28.94
(3,16)	1:B:21:C:O2	1:B:12:G:H22	11	28.94
(1,30)	1:B:22:A:H2	1:B:12:G:H1	18	28.94
(1,3)	1:B:12:G:H1	1:B:22:A:H2	18	28.94
(3,3)	1:B:12:G:H22	1:B:21:C:O2	1	28.92
(3,16)	1:B:21:C:O2	1:B:12:G:H22	1	28.92
(1,30)	1:B:22:A:H2	1:B:12:G:H1	15	28.92
(1,3)	1:B:12:G:H1	1:B:22:A:H2	15	28.92
(3,2)	1:B:12:G:O6	1:B:21:C:H42	13	28.91
(3,15)	1:B:21:C:H42	1:B:12:G:O6	13	28.91
(3,2)	1:B:12:G:O6	1:B:21:C:H42	5	28.87
(3,15)	1:B:21:C:H42	1:B:12:G:O6	5	28.87
(3,3)	1:B:12:G:H22	1:B:21:C:O2	2	28.85
(3,16)	1:B:21:C:O2	1:B:12:G:H22	2	28.85
(1,30)	1:B:22:A:H2	1:B:12:G:H1	5	28.84
(1,3)	1:B:12:G:H1	1:B:22:A:H2	5	28.84
(3,2)	1:B:12:G:O6	1:B:21:C:H42	1	28.83
(3,15)	1:B:21:C:H42	1:B:12:G:O6	1	28.83
(1,30)	1:B:22:A:H2	1:B:12:G:H1	17	28.81
(1,3)	1:B:12:G:H1	1:B:22:A:H2	17	28.81
(3,2)	1:B:12:G:O6	1:B:21:C:H42	16	28.8
(3,15)	1:B:21:C:H42	1:B:12:G:O6	16	28.8
(3,2)	1:B:12:G:O6	1:B:21:C:H42	18	28.79
(3,15)	1:B:21:C:H42	1:B:12:G:O6	18	28.79
(1,30)	1:B:22:A:H2	1:B:12:G:H1	11	28.79
(1,3)	1:B:12:G:H1	1:B:22:A:H2	11	28.79
(1,30)	1:B:22:A:H2	1:B:12:G:H1	19	28.78
(1,3)	1:B:12:G:H1	1:B:22:A:H2	19	28.78
(3,2)	1:B:12:G:O6	1:B:21:C:H42	12	28.76
(3,15)	1:B:21:C:H42	1:B:12:G:O6	12	28.76
(3,2)	1:B:12:G:O6	1:B:21:C:H42	7	28.75
(3,15)	1:B:21:C:H42	1:B:12:G:O6	7	28.75
(1,30)	1:B:22:A:H2	1:B:12:G:H1	8	28.75
(1,3)	1:B:12:G:H1	1:B:22:A:H2	8	28.75
(3,14)	1:B:21:C:N3	1:B:12:G:H1	15	28.73

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,1)	1:B:12:G:H1	1:B:21:C:N3	15	28.73
(1,30)	1:B:22:A:H2	1:B:12:G:H1	4	28.73
(1,3)	1:B:12:G:H1	1:B:22:A:H2	4	28.73
(1,30)	1:B:22:A:H2	1:B:12:G:H1	12	28.72
(1,3)	1:B:12:G:H1	1:B:22:A:H2	12	28.72
(3,2)	1:B:12:G:O6	1:B:21:C:H42	14	28.69
(3,15)	1:B:21:C:H42	1:B:12:G:O6	14	28.69
(3,14)	1:B:21:C:N3	1:B:12:G:H1	12	28.69
(3,1)	1:B:12:G:H1	1:B:21:C:N3	12	28.69
(3,14)	1:B:21:C:N3	1:B:12:G:H1	18	28.65
(3,14)	1:B:21:C:N3	1:B:12:G:H1	19	28.65
(3,1)	1:B:12:G:H1	1:B:21:C:N3	18	28.65
(3,1)	1:B:12:G:H1	1:B:21:C:N3	19	28.65
(3,14)	1:B:21:C:N3	1:B:12:G:H1	9	28.64
(3,14)	1:B:21:C:N3	1:B:12:G:H1	10	28.64
(3,1)	1:B:12:G:H1	1:B:21:C:N3	9	28.64
(3,1)	1:B:12:G:H1	1:B:21:C:N3	10	28.64
(3,14)	1:B:21:C:N3	1:B:12:G:H1	13	28.63
(3,1)	1:B:12:G:H1	1:B:21:C:N3	13	28.63
(3,2)	1:B:12:G:O6	1:B:21:C:H42	4	28.62
(3,15)	1:B:21:C:H42	1:B:12:G:O6	4	28.62
(3,14)	1:B:21:C:N3	1:B:12:G:H1	6	28.6
(3,1)	1:B:12:G:H1	1:B:21:C:N3	6	28.6
(1,30)	1:B:22:A:H2	1:B:12:G:H1	14	28.6
(1,3)	1:B:12:G:H1	1:B:22:A:H2	14	28.6
(3,14)	1:B:21:C:N3	1:B:12:G:H1	17	28.58
(3,1)	1:B:12:G:H1	1:B:21:C:N3	17	28.58
(3,14)	1:B:21:C:N3	1:B:12:G:H1	7	28.54
(3,14)	1:B:21:C:N3	1:B:12:G:H1	11	28.54
(3,14)	1:B:21:C:N3	1:B:12:G:H1	16	28.54
(3,1)	1:B:12:G:H1	1:B:21:C:N3	7	28.54
(3,1)	1:B:12:G:H1	1:B:21:C:N3	11	28.54
(3,1)	1:B:12:G:H1	1:B:21:C:N3	16	28.54
(3,14)	1:B:21:C:N3	1:B:12:G:H1	14	28.52
(3,1)	1:B:12:G:H1	1:B:21:C:N3	14	28.52
(1,30)	1:B:22:A:H2	1:B:12:G:H1	16	28.52
(1,3)	1:B:12:G:H1	1:B:22:A:H2	16	28.52
(1,30)	1:B:22:A:H2	1:B:12:G:H1	3	28.51
(1,3)	1:B:12:G:H1	1:B:22:A:H2	3	28.51
(3,14)	1:B:21:C:N3	1:B:12:G:H1	8	28.5
(3,1)	1:B:12:G:H1	1:B:21:C:N3	8	28.5
(3,14)	1:B:21:C:N3	1:B:12:G:H1	3	28.48

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,1)	1:B:12:G:H1	1:B:21:C:N3	3	28.48
(3,14)	1:B:21:C:N3	1:B:12:G:H1	5	28.47
(3,1)	1:B:12:G:H1	1:B:21:C:N3	5	28.47
(3,14)	1:B:21:C:N3	1:B:12:G:H1	4	28.44
(3,1)	1:B:12:G:H1	1:B:21:C:N3	4	28.44
(3,14)	1:B:21:C:N3	1:B:12:G:H1	2	28.41
(3,1)	1:B:12:G:H1	1:B:21:C:N3	2	28.41
(1,30)	1:B:22:A:H2	1:B:12:G:H1	2	28.38
(1,30)	1:B:22:A:H2	1:B:12:G:H1	10	28.38
(1,30)	1:B:22:A:H2	1:B:12:G:H1	13	28.38
(1,3)	1:B:12:G:H1	1:B:22:A:H2	2	28.38
(1,3)	1:B:12:G:H1	1:B:22:A:H2	10	28.38
(1,3)	1:B:12:G:H1	1:B:22:A:H2	13	28.38
(3,14)	1:B:21:C:N3	1:B:12:G:H1	1	28.31
(3,1)	1:B:12:G:H1	1:B:21:C:N3	1	28.31
(1,2)	1:B:12:G:H1	1:B:21:C:H42	11	28.01
(1,2)	1:B:12:G:H1	1:B:21:C:H42	15	28.01
(1,30)	1:B:22:A:H2	1:B:12:G:H1	1	27.96
(1,3)	1:B:12:G:H1	1:B:22:A:H2	1	27.96
(1,2)	1:B:12:G:H1	1:B:21:C:H42	9	27.88
(1,2)	1:B:12:G:H1	1:B:21:C:H42	10	27.81
(1,2)	1:B:12:G:H1	1:B:21:C:H42	12	27.78
(1,2)	1:B:12:G:H1	1:B:21:C:H42	3	27.77
(1,2)	1:B:12:G:H1	1:B:21:C:H42	5	27.75
(1,2)	1:B:12:G:H1	1:B:21:C:H42	2	27.74
(1,2)	1:B:12:G:H1	1:B:21:C:H42	6	27.74
(1,2)	1:B:12:G:H1	1:B:21:C:H42	17	27.72
(1,2)	1:B:12:G:H1	1:B:21:C:H42	8	27.71
(1,2)	1:B:12:G:H1	1:B:21:C:H42	19	27.69
(1,2)	1:B:12:G:H1	1:B:21:C:H42	13	27.67
(1,2)	1:B:12:G:H1	1:B:21:C:H42	18	27.66
(1,2)	1:B:12:G:H1	1:B:21:C:H42	4	27.56
(1,2)	1:B:12:G:H1	1:B:21:C:H42	7	27.52
(1,2)	1:B:12:G:H1	1:B:21:C:H42	16	27.52
(1,2)	1:B:12:G:H1	1:B:21:C:H42	1	27.42
(1,2)	1:B:12:G:H1	1:B:21:C:H42	14	27.29
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	13	24.41
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	13	24.41
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	12	24.36
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	12	24.36
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	10	24.16
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	10	24.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	14	24.14
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	14	24.14
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	16	24.05
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	16	24.05
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	7	24.04
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	7	24.04
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	3	24.01
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	3	24.01
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	18	23.97
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	18	23.97
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	4	23.95
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	4	23.95
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	19	23.94
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	19	23.94
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	6	23.91
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	6	23.91
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	1	23.85
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	1	23.85
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	17	23.82
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	17	23.82
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	8	23.79
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	8	23.79
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	15	23.73
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	15	23.73
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	2	23.7
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	2	23.7
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	11	23.67
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	11	23.67
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	9	23.65
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	9	23.65
(1,46)	1:B:13:A:H2	1:B:21:C:H1'	5	23.59
(1,108)	1:B:21:C:H1'	1:B:13:A:H2	5	23.59
(3,5)	1:B:13:A:H62	1:B:20:U:O4	15	23.4
(3,13)	1:B:20:U:O4	1:B:13:A:H62	15	23.4
(3,5)	1:B:13:A:H62	1:B:20:U:O4	17	23.3
(3,13)	1:B:20:U:O4	1:B:13:A:H62	17	23.3
(3,5)	1:B:13:A:H62	1:B:20:U:O4	19	23.17
(3,13)	1:B:20:U:O4	1:B:13:A:H62	19	23.17
(3,5)	1:B:13:A:H62	1:B:20:U:O4	9	23.12
(3,13)	1:B:20:U:O4	1:B:13:A:H62	9	23.12
(3,5)	1:B:13:A:H62	1:B:20:U:O4	11	23.09
(3,13)	1:B:20:U:O4	1:B:13:A:H62	11	23.09

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,5)	1:B:13:A:H62	1:B:20:U:O4	10	23.04
(3,13)	1:B:20:U:O4	1:B:13:A:H62	10	23.04
(3,5)	1:B:13:A:H62	1:B:20:U:O4	2	23.02
(3,13)	1:B:20:U:O4	1:B:13:A:H62	2	23.02
(3,5)	1:B:13:A:H62	1:B:20:U:O4	6	22.91
(3,13)	1:B:20:U:O4	1:B:13:A:H62	6	22.91
(3,5)	1:B:13:A:H62	1:B:20:U:O4	3	22.89
(3,5)	1:B:13:A:H62	1:B:20:U:O4	8	22.89
(3,13)	1:B:20:U:O4	1:B:13:A:H62	3	22.89
(3,13)	1:B:20:U:O4	1:B:13:A:H62	8	22.89
(3,5)	1:B:13:A:H62	1:B:20:U:O4	1	22.87
(3,13)	1:B:20:U:O4	1:B:13:A:H62	1	22.87
(3,5)	1:B:13:A:H62	1:B:20:U:O4	16	22.82
(3,13)	1:B:20:U:O4	1:B:13:A:H62	16	22.82
(3,5)	1:B:13:A:H62	1:B:20:U:O4	5	22.78
(3,13)	1:B:20:U:O4	1:B:13:A:H62	5	22.78
(3,5)	1:B:13:A:H62	1:B:20:U:O4	12	22.74
(3,13)	1:B:20:U:O4	1:B:13:A:H62	12	22.74
(3,5)	1:B:13:A:H62	1:B:20:U:O4	13	22.72
(3,13)	1:B:20:U:O4	1:B:13:A:H62	13	22.72
(3,5)	1:B:13:A:H62	1:B:20:U:O4	18	22.7
(3,13)	1:B:20:U:O4	1:B:13:A:H62	18	22.7
(3,5)	1:B:13:A:H62	1:B:20:U:O4	7	22.68
(3,13)	1:B:20:U:O4	1:B:13:A:H62	7	22.68
(3,5)	1:B:13:A:H62	1:B:20:U:O4	14	22.65
(3,13)	1:B:20:U:O4	1:B:13:A:H62	14	22.65
(3,5)	1:B:13:A:H62	1:B:20:U:O4	4	22.39
(3,13)	1:B:20:U:O4	1:B:13:A:H62	4	22.39
(3,4)	1:B:13:A:N1	1:B:20:U:H3	15	22.05
(3,12)	1:B:20:U:H3	1:B:13:A:N1	15	22.05
(3,4)	1:B:13:A:N1	1:B:20:U:H3	10	21.99
(3,12)	1:B:20:U:H3	1:B:13:A:N1	10	21.99
(3,4)	1:B:13:A:N1	1:B:20:U:H3	13	21.97
(3,12)	1:B:20:U:H3	1:B:13:A:N1	13	21.97
(3,4)	1:B:13:A:N1	1:B:20:U:H3	19	21.94
(3,12)	1:B:20:U:H3	1:B:13:A:N1	19	21.94
(3,4)	1:B:13:A:N1	1:B:20:U:H3	11	21.93
(3,12)	1:B:20:U:H3	1:B:13:A:N1	11	21.93
(3,4)	1:B:13:A:N1	1:B:20:U:H3	9	21.91
(3,12)	1:B:20:U:H3	1:B:13:A:N1	9	21.91
(3,4)	1:B:13:A:N1	1:B:20:U:H3	12	21.9
(3,12)	1:B:20:U:H3	1:B:13:A:N1	12	21.9

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,4)	1:B:13:A:N1	1:B:20:U:H3	3	21.88
(3,4)	1:B:13:A:N1	1:B:20:U:H3	17	21.88
(3,12)	1:B:20:U:H3	1:B:13:A:N1	3	21.88
(3,12)	1:B:20:U:H3	1:B:13:A:N1	17	21.88
(3,4)	1:B:13:A:N1	1:B:20:U:H3	6	21.86
(3,12)	1:B:20:U:H3	1:B:13:A:N1	6	21.86
(3,4)	1:B:13:A:N1	1:B:20:U:H3	2	21.83
(3,12)	1:B:20:U:H3	1:B:13:A:N1	2	21.83
(3,4)	1:B:13:A:N1	1:B:20:U:H3	8	21.8
(3,12)	1:B:20:U:H3	1:B:13:A:N1	8	21.8
(3,4)	1:B:13:A:N1	1:B:20:U:H3	16	21.77
(3,12)	1:B:20:U:H3	1:B:13:A:N1	16	21.77
(3,4)	1:B:13:A:N1	1:B:20:U:H3	5	21.73
(3,12)	1:B:20:U:H3	1:B:13:A:N1	5	21.73
(3,4)	1:B:13:A:N1	1:B:20:U:H3	7	21.69
(3,4)	1:B:13:A:N1	1:B:20:U:H3	18	21.69
(3,12)	1:B:20:U:H3	1:B:13:A:N1	7	21.69
(3,12)	1:B:20:U:H3	1:B:13:A:N1	18	21.69
(3,4)	1:B:13:A:N1	1:B:20:U:H3	1	21.68
(3,12)	1:B:20:U:H3	1:B:13:A:N1	1	21.68
(3,4)	1:B:13:A:N1	1:B:20:U:H3	14	21.62
(3,12)	1:B:20:U:H3	1:B:13:A:N1	14	21.62
(3,4)	1:B:13:A:N1	1:B:20:U:H3	4	21.59
(3,12)	1:B:20:U:H3	1:B:13:A:N1	4	21.59
(1,23)	1:B:20:U:H3	1:B:12:G:H1	15	21.13
(1,23)	1:B:20:U:H3	1:B:12:G:H1	10	21.06
(1,23)	1:B:20:U:H3	1:B:12:G:H1	13	21.03
(1,23)	1:B:20:U:H3	1:B:12:G:H1	19	21.03
(1,23)	1:B:20:U:H3	1:B:12:G:H1	9	20.95
(1,23)	1:B:20:U:H3	1:B:12:G:H1	12	20.95
(1,23)	1:B:20:U:H3	1:B:12:G:H1	11	20.94
(1,23)	1:B:20:U:H3	1:B:12:G:H1	17	20.93
(1,23)	1:B:20:U:H3	1:B:12:G:H1	6	20.9
(1,23)	1:B:20:U:H3	1:B:12:G:H1	8	20.89
(1,23)	1:B:20:U:H3	1:B:12:G:H1	2	20.87
(1,23)	1:B:20:U:H3	1:B:12:G:H1	3	20.87
(1,23)	1:B:20:U:H3	1:B:12:G:H1	18	20.86
(1,23)	1:B:20:U:H3	1:B:12:G:H1	16	20.83
(1,23)	1:B:20:U:H3	1:B:12:G:H1	5	20.8
(1,23)	1:B:20:U:H3	1:B:12:G:H1	1	20.75
(1,23)	1:B:20:U:H3	1:B:12:G:H1	7	20.7
(1,23)	1:B:20:U:H3	1:B:12:G:H1	4	20.68

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,23)	1:B:20:U:H3	1:B:12:G:H1	14	20.64
(1,24)	1:B:20:U:H3	1:B:13:A:H2	12	20.51
(1,24)	1:B:20:U:H3	1:B:13:A:H2	3	20.46
(1,24)	1:B:20:U:H3	1:B:13:A:H2	13	20.45
(1,24)	1:B:20:U:H3	1:B:13:A:H2	11	20.42
(1,1)	1:B:12:G:H1	1:B:20:U:H3	15	20.41
(1,24)	1:B:20:U:H3	1:B:13:A:H2	10	20.4
(1,24)	1:B:20:U:H3	1:B:13:A:H2	6	20.35
(1,1)	1:B:12:G:H1	1:B:20:U:H3	10	20.34
(1,24)	1:B:20:U:H3	1:B:13:A:H2	9	20.33
(1,24)	1:B:20:U:H3	1:B:13:A:H2	15	20.33
(1,1)	1:B:12:G:H1	1:B:20:U:H3	13	20.31
(1,1)	1:B:12:G:H1	1:B:20:U:H3	19	20.31
(1,24)	1:B:20:U:H3	1:B:13:A:H2	5	20.27
(1,24)	1:B:20:U:H3	1:B:13:A:H2	2	20.25
(1,24)	1:B:20:U:H3	1:B:13:A:H2	4	20.23
(1,24)	1:B:20:U:H3	1:B:13:A:H2	16	20.23
(1,1)	1:B:12:G:H1	1:B:20:U:H3	9	20.23
(1,1)	1:B:12:G:H1	1:B:20:U:H3	12	20.23
(1,1)	1:B:12:G:H1	1:B:20:U:H3	11	20.22
(1,24)	1:B:20:U:H3	1:B:13:A:H2	8	20.21
(1,1)	1:B:12:G:H1	1:B:20:U:H3	17	20.21
(1,24)	1:B:20:U:H3	1:B:13:A:H2	19	20.2
(1,24)	1:B:20:U:H3	1:B:13:A:H2	18	20.19
(1,1)	1:B:12:G:H1	1:B:20:U:H3	6	20.18
(1,24)	1:B:20:U:H3	1:B:13:A:H2	7	20.17
(1,1)	1:B:12:G:H1	1:B:20:U:H3	8	20.17
(1,1)	1:B:12:G:H1	1:B:20:U:H3	2	20.15
(1,1)	1:B:12:G:H1	1:B:20:U:H3	3	20.15
(1,1)	1:B:12:G:H1	1:B:20:U:H3	18	20.14
(1,1)	1:B:12:G:H1	1:B:20:U:H3	16	20.11
(1,24)	1:B:20:U:H3	1:B:13:A:H2	17	20.1
(1,1)	1:B:12:G:H1	1:B:20:U:H3	5	20.08
(1,24)	1:B:20:U:H3	1:B:13:A:H2	14	20.03
(1,1)	1:B:12:G:H1	1:B:20:U:H3	1	20.03
(1,24)	1:B:20:U:H3	1:B:13:A:H2	1	19.99
(1,1)	1:B:12:G:H1	1:B:20:U:H3	7	19.98
(1,1)	1:B:12:G:H1	1:B:20:U:H3	4	19.96
(1,1)	1:B:12:G:H1	1:B:20:U:H3	14	19.92
(1,9)	1:B:14:C:H42	1:B:20:U:H3	15	17.06
(1,25)	1:B:20:U:H3	1:B:14:C:H42	15	17.06
(1,9)	1:B:14:C:H42	1:B:20:U:H3	19	16.94

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,25)	1:B:20:U:H3	1:B:14:C:H42	19	16.94
(3,7)	1:B:14:C:H42	1:B:19:G:O6	15	16.92
(3,10)	1:B:19:G:O6	1:B:14:C:H42	15	16.92
(3,7)	1:B:14:C:H42	1:B:19:G:O6	9	16.88
(3,10)	1:B:19:G:O6	1:B:14:C:H42	9	16.88
(1,9)	1:B:14:C:H42	1:B:20:U:H3	9	16.88
(1,25)	1:B:20:U:H3	1:B:14:C:H42	9	16.88
(3,7)	1:B:14:C:H42	1:B:19:G:O6	11	16.82
(3,10)	1:B:19:G:O6	1:B:14:C:H42	11	16.82
(1,9)	1:B:14:C:H42	1:B:20:U:H3	6	16.82
(1,25)	1:B:20:U:H3	1:B:14:C:H42	6	16.82
(1,9)	1:B:14:C:H42	1:B:20:U:H3	10	16.78
(1,9)	1:B:14:C:H42	1:B:20:U:H3	17	16.78
(1,25)	1:B:20:U:H3	1:B:14:C:H42	10	16.78
(1,25)	1:B:20:U:H3	1:B:14:C:H42	17	16.78
(1,9)	1:B:14:C:H42	1:B:20:U:H3	11	16.76
(1,25)	1:B:20:U:H3	1:B:14:C:H42	11	16.76
(3,7)	1:B:14:C:H42	1:B:19:G:O6	6	16.75
(3,10)	1:B:19:G:O6	1:B:14:C:H42	6	16.75
(3,7)	1:B:14:C:H42	1:B:19:G:O6	17	16.73
(3,10)	1:B:19:G:O6	1:B:14:C:H42	17	16.73
(3,7)	1:B:14:C:H42	1:B:19:G:O6	19	16.68
(3,10)	1:B:19:G:O6	1:B:14:C:H42	19	16.68
(1,9)	1:B:14:C:H42	1:B:20:U:H3	1	16.68
(1,25)	1:B:20:U:H3	1:B:14:C:H42	1	16.68
(1,9)	1:B:14:C:H42	1:B:20:U:H3	13	16.67
(1,25)	1:B:20:U:H3	1:B:14:C:H42	13	16.67
(1,9)	1:B:14:C:H42	1:B:20:U:H3	16	16.65
(1,25)	1:B:20:U:H3	1:B:14:C:H42	16	16.65
(1,9)	1:B:14:C:H42	1:B:20:U:H3	2	16.58
(1,25)	1:B:20:U:H3	1:B:14:C:H42	2	16.58
(3,7)	1:B:14:C:H42	1:B:19:G:O6	10	16.57
(3,10)	1:B:19:G:O6	1:B:14:C:H42	10	16.57
(3,7)	1:B:14:C:H42	1:B:19:G:O6	16	16.55
(3,10)	1:B:19:G:O6	1:B:14:C:H42	16	16.55
(1,9)	1:B:14:C:H42	1:B:20:U:H3	8	16.52
(1,9)	1:B:14:C:H42	1:B:20:U:H3	14	16.52
(1,25)	1:B:20:U:H3	1:B:14:C:H42	8	16.52
(1,25)	1:B:20:U:H3	1:B:14:C:H42	14	16.52
(1,9)	1:B:14:C:H42	1:B:20:U:H3	5	16.51
(1,25)	1:B:20:U:H3	1:B:14:C:H42	5	16.51
(3,7)	1:B:14:C:H42	1:B:19:G:O6	2	16.49

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,7)	1:B:14:C:H42	1:B:19:G:O6	5	16.49
(3,10)	1:B:19:G:O6	1:B:14:C:H42	2	16.49
(3,10)	1:B:19:G:O6	1:B:14:C:H42	5	16.49
(1,9)	1:B:14:C:H42	1:B:20:U:H3	12	16.49
(1,25)	1:B:20:U:H3	1:B:14:C:H42	12	16.49
(1,9)	1:B:14:C:H42	1:B:20:U:H3	3	16.47
(1,25)	1:B:20:U:H3	1:B:14:C:H42	3	16.47
(3,7)	1:B:14:C:H42	1:B:19:G:O6	3	16.45
(3,10)	1:B:19:G:O6	1:B:14:C:H42	3	16.45
(3,7)	1:B:14:C:H42	1:B:19:G:O6	8	16.37
(3,10)	1:B:19:G:O6	1:B:14:C:H42	8	16.37
(1,9)	1:B:14:C:H42	1:B:20:U:H3	18	16.37
(1,25)	1:B:20:U:H3	1:B:14:C:H42	18	16.37
(3,7)	1:B:14:C:H42	1:B:19:G:O6	1	16.36
(3,10)	1:B:19:G:O6	1:B:14:C:H42	1	16.36
(3,7)	1:B:14:C:H42	1:B:19:G:O6	12	16.33
(3,10)	1:B:19:G:O6	1:B:14:C:H42	12	16.33
(3,7)	1:B:14:C:H42	1:B:19:G:O6	14	16.32
(3,10)	1:B:19:G:O6	1:B:14:C:H42	14	16.32
(1,9)	1:B:14:C:H42	1:B:20:U:H3	7	16.29
(1,25)	1:B:20:U:H3	1:B:14:C:H42	7	16.29
(3,7)	1:B:14:C:H42	1:B:19:G:O6	13	16.24
(3,7)	1:B:14:C:H42	1:B:19:G:O6	18	16.24
(3,10)	1:B:19:G:O6	1:B:14:C:H42	13	16.24
(3,10)	1:B:19:G:O6	1:B:14:C:H42	18	16.24
(3,7)	1:B:14:C:H42	1:B:19:G:O6	7	16.19
(3,10)	1:B:19:G:O6	1:B:14:C:H42	7	16.19
(1,9)	1:B:14:C:H42	1:B:20:U:H3	4	16.08
(1,25)	1:B:20:U:H3	1:B:14:C:H42	4	16.08
(3,7)	1:B:14:C:H42	1:B:19:G:O6	4	16.04
(3,10)	1:B:19:G:O6	1:B:14:C:H42	4	16.04
(1,19)	1:B:19:G:H1	1:B:14:C:H42	15	15.3
(3,9)	1:B:19:G:H1	1:B:14:C:N3	15	15.22
(3,6)	1:B:14:C:N3	1:B:19:G:H1	15	15.22
(3,9)	1:B:19:G:H1	1:B:14:C:N3	10	15.21
(3,6)	1:B:14:C:N3	1:B:19:G:H1	10	15.21
(1,19)	1:B:19:G:H1	1:B:14:C:H42	19	15.21
(3,9)	1:B:19:G:H1	1:B:14:C:N3	19	15.2
(3,6)	1:B:14:C:N3	1:B:19:G:H1	19	15.2
(1,19)	1:B:19:G:H1	1:B:14:C:H42	9	15.14
(3,9)	1:B:19:G:H1	1:B:14:C:N3	11	15.13
(3,9)	1:B:19:G:H1	1:B:14:C:N3	13	15.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,6)	1:B:14:C:N3	1:B:19:G:H1	11	15.13
(3,6)	1:B:14:C:N3	1:B:19:G:H1	13	15.13
(3,9)	1:B:19:G:H1	1:B:14:C:N3	9	15.11
(3,9)	1:B:19:G:H1	1:B:14:C:N3	17	15.11
(3,6)	1:B:14:C:N3	1:B:19:G:H1	9	15.11
(3,6)	1:B:14:C:N3	1:B:19:G:H1	17	15.11
(3,8)	1:B:14:C:O2	1:B:19:G:H22	13	15.08
(3,11)	1:B:19:G:H22	1:B:14:C:O2	13	15.08
(3,9)	1:B:19:G:H1	1:B:14:C:N3	3	15.05
(3,6)	1:B:14:C:N3	1:B:19:G:H1	3	15.05
(1,19)	1:B:19:G:H1	1:B:14:C:H42	10	15.05
(1,19)	1:B:19:G:H1	1:B:14:C:H42	11	15.04
(1,19)	1:B:19:G:H1	1:B:14:C:H42	6	15.03
(1,19)	1:B:19:G:H1	1:B:14:C:H42	17	15.03
(3,9)	1:B:19:G:H1	1:B:14:C:N3	2	15.02
(3,6)	1:B:14:C:N3	1:B:19:G:H1	2	15.02
(3,9)	1:B:19:G:H1	1:B:14:C:N3	1	14.98
(3,9)	1:B:19:G:H1	1:B:14:C:N3	6	14.98
(3,6)	1:B:14:C:N3	1:B:19:G:H1	1	14.98
(3,6)	1:B:14:C:N3	1:B:19:G:H1	6	14.98
(3,9)	1:B:19:G:H1	1:B:14:C:N3	16	14.97
(3,6)	1:B:14:C:N3	1:B:19:G:H1	16	14.97
(1,19)	1:B:19:G:H1	1:B:14:C:H42	1	14.97
(1,19)	1:B:19:G:H1	1:B:14:C:H42	13	14.97
(3,8)	1:B:14:C:O2	1:B:19:G:H22	10	14.96
(3,11)	1:B:19:G:H22	1:B:14:C:O2	10	14.96
(3,9)	1:B:19:G:H1	1:B:14:C:N3	8	14.95
(3,9)	1:B:19:G:H1	1:B:14:C:N3	12	14.95
(3,6)	1:B:14:C:N3	1:B:19:G:H1	8	14.95
(3,6)	1:B:14:C:N3	1:B:19:G:H1	12	14.95
(1,19)	1:B:19:G:H1	1:B:14:C:H42	16	14.95
(1,19)	1:B:19:G:H1	1:B:14:C:H42	2	14.92
(3,8)	1:B:14:C:O2	1:B:19:G:H22	1	14.88
(3,11)	1:B:19:G:H22	1:B:14:C:O2	1	14.88
(3,9)	1:B:19:G:H1	1:B:14:C:N3	5	14.87
(3,9)	1:B:19:G:H1	1:B:14:C:N3	7	14.87
(3,6)	1:B:14:C:N3	1:B:19:G:H1	5	14.87
(3,6)	1:B:14:C:N3	1:B:19:G:H1	7	14.87
(3,9)	1:B:19:G:H1	1:B:14:C:N3	18	14.84
(3,6)	1:B:14:C:N3	1:B:19:G:H1	18	14.84
(3,9)	1:B:19:G:H1	1:B:14:C:N3	14	14.83
(3,8)	1:B:14:C:O2	1:B:19:G:H22	19	14.83

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(3,6)	1:B:14:C:N3	1:B:19:G:H1	14	14.83
(3,11)	1:B:19:G:H22	1:B:14:C:O2	19	14.83
(1,19)	1:B:19:G:H1	1:B:14:C:H42	8	14.83
(3,9)	1:B:19:G:H1	1:B:14:C:N3	4	14.79
(3,6)	1:B:14:C:N3	1:B:19:G:H1	4	14.79
(1,19)	1:B:19:G:H1	1:B:14:C:H42	14	14.79
(1,19)	1:B:19:G:H1	1:B:14:C:H42	3	14.76
(1,19)	1:B:19:G:H1	1:B:14:C:H42	5	14.71
(3,8)	1:B:14:C:O2	1:B:19:G:H22	12	14.7
(3,8)	1:B:14:C:O2	1:B:19:G:H22	16	14.7
(3,8)	1:B:14:C:O2	1:B:19:G:H22	17	14.7
(3,11)	1:B:19:G:H22	1:B:14:C:O2	12	14.7
(3,11)	1:B:19:G:H22	1:B:14:C:O2	16	14.7
(3,11)	1:B:19:G:H22	1:B:14:C:O2	17	14.7
(3,8)	1:B:14:C:O2	1:B:19:G:H22	3	14.69
(3,11)	1:B:19:G:H22	1:B:14:C:O2	3	14.69
(3,8)	1:B:14:C:O2	1:B:19:G:H22	8	14.67
(3,11)	1:B:19:G:H22	1:B:14:C:O2	8	14.67
(1,19)	1:B:19:G:H1	1:B:14:C:H42	12	14.67
(3,8)	1:B:14:C:O2	1:B:19:G:H22	2	14.65
(3,11)	1:B:19:G:H22	1:B:14:C:O2	2	14.65
(1,19)	1:B:19:G:H1	1:B:14:C:H42	18	14.62
(3,8)	1:B:14:C:O2	1:B:19:G:H22	15	14.58
(3,11)	1:B:19:G:H22	1:B:14:C:O2	15	14.58
(3,8)	1:B:14:C:O2	1:B:19:G:H22	4	14.56
(3,11)	1:B:19:G:H22	1:B:14:C:O2	4	14.56
(3,8)	1:B:14:C:O2	1:B:19:G:H22	11	14.55
(3,11)	1:B:19:G:H22	1:B:14:C:O2	11	14.55
(3,8)	1:B:14:C:O2	1:B:19:G:H22	7	14.51
(3,11)	1:B:19:G:H22	1:B:14:C:O2	7	14.51
(3,8)	1:B:14:C:O2	1:B:19:G:H22	14	14.47
(3,11)	1:B:19:G:H22	1:B:14:C:O2	14	14.47
(1,19)	1:B:19:G:H1	1:B:14:C:H42	7	14.47
(3,8)	1:B:14:C:O2	1:B:19:G:H22	18	14.46
(3,11)	1:B:19:G:H22	1:B:14:C:O2	18	14.46
(3,8)	1:B:14:C:O2	1:B:19:G:H22	9	14.45
(3,11)	1:B:19:G:H22	1:B:14:C:O2	9	14.45
(3,8)	1:B:14:C:O2	1:B:19:G:H22	5	14.43
(3,11)	1:B:19:G:H22	1:B:14:C:O2	5	14.43
(3,8)	1:B:14:C:O2	1:B:19:G:H22	6	14.39
(3,11)	1:B:19:G:H22	1:B:14:C:O2	6	14.39
(1,19)	1:B:19:G:H1	1:B:14:C:H42	4	14.27

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,20)	1:B:19:G:H1	1:B:15:G:H8	17	7.83
(1,13)	1:B:15:G:H8	1:B:19:G:H1	17	7.83
(1,20)	1:B:19:G:H1	1:B:15:G:H8	3	7.7
(1,13)	1:B:15:G:H8	1:B:19:G:H1	3	7.7
(1,20)	1:B:19:G:H1	1:B:15:G:H8	7	7.65
(1,13)	1:B:15:G:H8	1:B:19:G:H1	7	7.65
(1,20)	1:B:19:G:H1	1:B:15:G:H8	10	7.61
(1,13)	1:B:15:G:H8	1:B:19:G:H1	10	7.61
(1,20)	1:B:19:G:H1	1:B:15:G:H8	13	7.57
(1,13)	1:B:15:G:H8	1:B:19:G:H1	13	7.57
(1,20)	1:B:19:G:H1	1:B:15:G:H8	2	7.46
(1,13)	1:B:15:G:H8	1:B:19:G:H1	2	7.46
(1,20)	1:B:19:G:H1	1:B:15:G:H8	1	7.45
(1,20)	1:B:19:G:H1	1:B:15:G:H8	14	7.45
(1,13)	1:B:15:G:H8	1:B:19:G:H1	1	7.45
(1,13)	1:B:15:G:H8	1:B:19:G:H1	14	7.45
(1,20)	1:B:19:G:H1	1:B:15:G:H8	19	7.43
(1,13)	1:B:15:G:H8	1:B:19:G:H1	19	7.43
(1,20)	1:B:19:G:H1	1:B:15:G:H8	8	7.38
(1,13)	1:B:15:G:H8	1:B:19:G:H1	8	7.38
(1,20)	1:B:19:G:H1	1:B:15:G:H8	6	7.37
(1,13)	1:B:15:G:H8	1:B:19:G:H1	6	7.37
(1,20)	1:B:19:G:H1	1:B:15:G:H8	15	7.35
(1,13)	1:B:15:G:H8	1:B:19:G:H1	15	7.35
(1,20)	1:B:19:G:H1	1:B:15:G:H8	16	7.34
(1,13)	1:B:15:G:H8	1:B:19:G:H1	16	7.34
(1,20)	1:B:19:G:H1	1:B:15:G:H8	5	7.33
(1,13)	1:B:15:G:H8	1:B:19:G:H1	5	7.33
(1,20)	1:B:19:G:H1	1:B:15:G:H8	9	7.31
(1,13)	1:B:15:G:H8	1:B:19:G:H1	9	7.31
(1,20)	1:B:19:G:H1	1:B:15:G:H8	11	7.27
(1,13)	1:B:15:G:H8	1:B:19:G:H1	11	7.27
(1,20)	1:B:19:G:H1	1:B:15:G:H8	12	7.22
(1,20)	1:B:19:G:H1	1:B:15:G:H8	18	7.22
(1,13)	1:B:15:G:H8	1:B:19:G:H1	12	7.22
(1,13)	1:B:15:G:H8	1:B:19:G:H1	18	7.22
(1,20)	1:B:19:G:H1	1:B:15:G:H8	4	7.06
(1,13)	1:B:15:G:H8	1:B:19:G:H1	4	7.06
(1,34)	1:B:17:G:H1	1:B:15:G:H8	17	6.27
(1,32)	1:B:15:G:H8	1:B:17:G:H1	17	6.27
(1,34)	1:B:17:G:H1	1:B:15:G:H8	10	6.25
(1,32)	1:B:15:G:H8	1:B:17:G:H1	10	6.25

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,34)	1:B:17:G:H1	1:B:15:G:H8	13	6.18
(1,32)	1:B:15:G:H8	1:B:17:G:H1	13	6.18
(1,34)	1:B:17:G:H1	1:B:15:G:H8	3	6.16
(1,34)	1:B:17:G:H1	1:B:15:G:H8	16	6.16
(1,32)	1:B:15:G:H8	1:B:17:G:H1	3	6.16
(1,32)	1:B:15:G:H8	1:B:17:G:H1	16	6.16
(1,34)	1:B:17:G:H1	1:B:15:G:H8	1	6.13
(1,32)	1:B:15:G:H8	1:B:17:G:H1	1	6.13
(1,34)	1:B:17:G:H1	1:B:15:G:H8	2	6.12
(1,34)	1:B:17:G:H1	1:B:15:G:H8	7	6.12
(1,32)	1:B:15:G:H8	1:B:17:G:H1	2	6.12
(1,32)	1:B:15:G:H8	1:B:17:G:H1	7	6.12
(1,34)	1:B:17:G:H1	1:B:15:G:H8	14	6.11
(1,32)	1:B:15:G:H8	1:B:17:G:H1	14	6.11
(1,34)	1:B:17:G:H1	1:B:15:G:H8	19	6.06
(1,32)	1:B:15:G:H8	1:B:17:G:H1	19	6.06
(1,34)	1:B:17:G:H1	1:B:15:G:H8	5	6.02
(1,32)	1:B:15:G:H8	1:B:17:G:H1	5	6.02
(1,34)	1:B:17:G:H1	1:B:15:G:H8	6	5.93
(1,32)	1:B:15:G:H8	1:B:17:G:H1	6	5.93
(1,34)	1:B:17:G:H1	1:B:15:G:H8	15	5.92
(1,32)	1:B:15:G:H8	1:B:17:G:H1	15	5.92
(1,34)	1:B:17:G:H1	1:B:15:G:H8	11	5.88
(1,32)	1:B:15:G:H8	1:B:17:G:H1	11	5.88
(1,34)	1:B:17:G:H1	1:B:15:G:H8	12	5.87
(1,32)	1:B:15:G:H8	1:B:17:G:H1	12	5.87
(1,34)	1:B:17:G:H1	1:B:15:G:H8	8	5.86
(1,32)	1:B:15:G:H8	1:B:17:G:H1	8	5.86
(1,34)	1:B:17:G:H1	1:B:15:G:H8	9	5.84
(1,32)	1:B:15:G:H8	1:B:17:G:H1	9	5.84
(1,34)	1:B:17:G:H1	1:B:15:G:H8	4	5.8
(1,32)	1:B:15:G:H8	1:B:17:G:H1	4	5.8
(1,34)	1:B:17:G:H1	1:B:15:G:H8	18	5.74
(1,32)	1:B:15:G:H8	1:B:17:G:H1	18	5.74
(1,33)	1:B:16:A:H2	1:B:15:G:H1	16	4.24
(1,31)	1:B:15:G:H1	1:B:16:A:H2	16	4.24
(1,33)	1:B:16:A:H2	1:B:15:G:H1	9	4.14
(1,31)	1:B:15:G:H1	1:B:16:A:H2	9	4.14
(1,33)	1:B:16:A:H2	1:B:15:G:H1	5	4.13
(1,31)	1:B:15:G:H1	1:B:16:A:H2	5	4.13
(1,33)	1:B:16:A:H2	1:B:15:G:H1	19	4.12
(1,31)	1:B:15:G:H1	1:B:16:A:H2	19	4.12

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,33)	1:B:16:A:H2	1:B:15:G:H1	6	4.06
(1,31)	1:B:15:G:H1	1:B:16:A:H2	6	4.06
(1,33)	1:B:16:A:H2	1:B:15:G:H1	12	3.95
(1,31)	1:B:15:G:H1	1:B:16:A:H2	12	3.95
(1,33)	1:B:16:A:H2	1:B:15:G:H1	7	3.9
(1,31)	1:B:15:G:H1	1:B:16:A:H2	7	3.9
(1,33)	1:B:16:A:H2	1:B:15:G:H1	4	3.77
(1,33)	1:B:16:A:H2	1:B:15:G:H1	10	3.77
(1,31)	1:B:15:G:H1	1:B:16:A:H2	4	3.77
(1,31)	1:B:15:G:H1	1:B:16:A:H2	10	3.77
(1,33)	1:B:16:A:H2	1:B:15:G:H1	15	3.72
(1,31)	1:B:15:G:H1	1:B:16:A:H2	15	3.72
(1,33)	1:B:16:A:H2	1:B:15:G:H1	18	3.71
(1,31)	1:B:15:G:H1	1:B:16:A:H2	18	3.71
(1,33)	1:B:16:A:H2	1:B:15:G:H1	14	3.61
(1,31)	1:B:15:G:H1	1:B:16:A:H2	14	3.61
(1,33)	1:B:16:A:H2	1:B:15:G:H1	1	3.59
(1,31)	1:B:15:G:H1	1:B:16:A:H2	1	3.59
(1,33)	1:B:16:A:H2	1:B:15:G:H1	8	3.57
(1,33)	1:B:16:A:H2	1:B:15:G:H1	11	3.57
(1,31)	1:B:15:G:H1	1:B:16:A:H2	8	3.57
(1,31)	1:B:15:G:H1	1:B:16:A:H2	11	3.57
(1,33)	1:B:16:A:H2	1:B:15:G:H1	17	3.4
(1,31)	1:B:15:G:H1	1:B:16:A:H2	17	3.4
(1,33)	1:B:16:A:H2	1:B:15:G:H1	2	3.39
(1,31)	1:B:15:G:H1	1:B:16:A:H2	2	3.39
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	13	3.36
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	13	3.36
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	10	3.33
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	10	3.33
(1,33)	1:B:16:A:H2	1:B:15:G:H1	13	3.32
(1,31)	1:B:15:G:H1	1:B:16:A:H2	13	3.32
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	17	3.32
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	17	3.32
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	1	3.31
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	1	3.31
(1,33)	1:B:16:A:H2	1:B:15:G:H1	3	3.27
(1,31)	1:B:15:G:H1	1:B:16:A:H2	3	3.27
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	16	3.27
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	16	3.27
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	14	3.25
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	14	3.25

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	2	3.23
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	3	3.23
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	2	3.23
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	3	3.23
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	7	3.15
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	7	3.15
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	19	3.12
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	19	3.12
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	5	3.08
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	8	3.08
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	5	3.08
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	8	3.08
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	12	3.07
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	12	3.07
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	6	3.04
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	6	3.04
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	9	2.99
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	15	2.99
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	9	2.99
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	15	2.99
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	18	2.95
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	18	2.95
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	4	2.93
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	4	2.93
(1,16)	1:B:17:G:H1	1:B:15:G:H1'	11	2.92
(1,12)	1:B:15:G:H1'	1:B:17:G:H1	11	2.92
(1,141)	1:B:16:A:N7	1:B:16:A:H61	4	1.1
(1,140)	1:B:13:A:N7	1:B:13:A:H61	2	1.1
(1,142)	1:B:22:A:N7	1:B:22:A:H61	5	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	6	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	8	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	10	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	11	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	12	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	13	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	14	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	17	1.09
(1,141)	1:B:16:A:N7	1:B:16:A:H61	1	1.09
(1,141)	1:B:16:A:N7	1:B:16:A:H61	7	1.09
(1,141)	1:B:16:A:N7	1:B:16:A:H61	8	1.09
(1,141)	1:B:16:A:N7	1:B:16:A:H61	9	1.09
(1,141)	1:B:16:A:N7	1:B:16:A:H61	12	1.09

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,141)	1:B:16:A:N7	1:B:16:A:H61	18	1.09
(1,140)	1:B:13:A:N7	1:B:13:A:H61	1	1.09
(1,140)	1:B:13:A:N7	1:B:13:A:H61	5	1.09
(1,140)	1:B:13:A:N7	1:B:13:A:H61	8	1.09
(1,140)	1:B:13:A:N7	1:B:13:A:H61	12	1.09
(1,140)	1:B:13:A:N7	1:B:13:A:H61	17	1.09
(1,142)	1:B:22:A:N7	1:B:22:A:H61	1	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	2	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	3	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	4	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	7	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	9	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	15	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	16	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	18	1.08
(1,142)	1:B:22:A:N7	1:B:22:A:H61	19	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	3	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	5	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	6	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	11	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	13	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	14	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	15	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	16	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	17	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	19	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	3	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	4	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	6	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	7	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	9	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	10	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	13	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	15	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	16	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	18	1.08
(1,140)	1:B:13:A:N7	1:B:13:A:H61	19	1.08
(1,141)	1:B:16:A:N7	1:B:16:A:H61	2	1.07
(1,141)	1:B:16:A:N7	1:B:16:A:H61	10	1.07
(1,140)	1:B:13:A:N7	1:B:13:A:H61	11	1.07
(1,140)	1:B:13:A:N7	1:B:13:A:H61	14	1.07
(1,139)	1:B:21:C:H5	1:B:21:C:H41	3	1.01

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,139)	1:B:21:C:H5	1:B:21:C:H41	6	1.01
(1,139)	1:B:21:C:H5	1:B:21:C:H41	14	1.01
(1,139)	1:B:21:C:H5	1:B:21:C:H41	17	1.01
(1,139)	1:B:21:C:H5	1:B:21:C:H41	4	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	5	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	7	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	8	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	9	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	10	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	11	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	12	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	15	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	16	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	18	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	19	1.0
(1,139)	1:B:21:C:H5	1:B:21:C:H41	1	0.99
(1,139)	1:B:21:C:H5	1:B:21:C:H41	2	0.99
(1,139)	1:B:21:C:H5	1:B:21:C:H41	13	0.99
(1,138)	1:B:14:C:H5	1:B:14:C:H41	2	0.99
(1,138)	1:B:14:C:H5	1:B:14:C:H41	3	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	5	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	6	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	7	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	8	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	11	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	12	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	13	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	14	0.98
(1,138)	1:B:14:C:H5	1:B:14:C:H41	1	0.97
(1,138)	1:B:14:C:H5	1:B:14:C:H41	4	0.97
(1,138)	1:B:14:C:H5	1:B:14:C:H41	9	0.97
(1,138)	1:B:14:C:H5	1:B:14:C:H41	10	0.97
(1,138)	1:B:14:C:H5	1:B:14:C:H41	15	0.97
(1,138)	1:B:14:C:H5	1:B:14:C:H41	16	0.97
(1,138)	1:B:14:C:H5	1:B:14:C:H41	17	0.97
(1,138)	1:B:14:C:H5	1:B:14:C:H41	18	0.97
(1,138)	1:B:14:C:H5	1:B:14:C:H41	19	0.97
(1,136)	1:B:17:G:H1	1:B:17:G:H22	2	0.95
(1,136)	1:B:17:G:H1	1:B:17:G:H22	5	0.95
(1,135)	1:B:15:G:H1	1:B:15:G:H22	2	0.95
(1,135)	1:B:15:G:H1	1:B:15:G:H22	6	0.95
(1,135)	1:B:15:G:H1	1:B:15:G:H22	10	0.95

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,135)	1:B:15:G:H1	1:B:15:G:H22	11	0.95
(1,135)	1:B:15:G:H1	1:B:15:G:H22	12	0.95
(1,135)	1:B:15:G:H1	1:B:15:G:H22	18	0.95
(1,135)	1:B:15:G:H1	1:B:15:G:H22	19	0.95
(1,135)	1:B:15:G:H1	1:B:15:G:H22	1	0.94
(1,135)	1:B:15:G:H1	1:B:15:G:H22	5	0.94
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	10	0.54
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	10	0.54
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	1	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	1	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	2	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	2	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	3	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	3	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	5	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	5	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	6	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	6	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	7	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	7	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	8	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	8	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	9	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	9	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	11	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	11	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	12	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	12	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	15	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	15	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	16	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	16	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	17	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	17	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	18	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	18	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	19	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	19	0.53
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	4	0.52
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	4	0.52
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	13	0.52
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	13	0.52

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	14	0.52
(1,44)	1:B:13:A:H1'	1:B:13:A:H2	14	0.52
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	4	0.28
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	4	0.28
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	10	0.28
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	10	0.28
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	7	0.27
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	7	0.27
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	12	0.27
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	12	0.27
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	2	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	2	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	8	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	8	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	13	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	13	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	14	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	14	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	16	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	16	0.26
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	17	0.25
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	17	0.25
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	19	0.25
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	19	0.25
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	1	0.24
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	1	0.24
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	10	0.24
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	10	0.24
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	15	0.24
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	15	0.24
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	3	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	3	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	4	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	4	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	6	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	6	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	11	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	11	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	18	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	18	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	19	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	19	0.23

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	11	0.23
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	11	0.23
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	18	0.23
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	18	0.23
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	5	0.22
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	5	0.22
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	12	0.22
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	12	0.22
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	5	0.22
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	5	0.22
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	8	0.22
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	8	0.22
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	17	0.22
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	17	0.22
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	7	0.21
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	7	0.21
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	10	0.21
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	10	0.21
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	3	0.21
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	3	0.21
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	12	0.2
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	12	0.2
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	9	0.19
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	9	0.19
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	18	0.19
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	18	0.19
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	9	0.19
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	9	0.19
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	13	0.18
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	13	0.18
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	14	0.18
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	14	0.18
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	16	0.18
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	16	0.18
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	1	0.18
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	1	0.18
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	6	0.18
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	6	0.18
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	13	0.18
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	13	0.18
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	4	0.18
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	4	0.18

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	16	0.18
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	16	0.18
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	7	0.17
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	7	0.17
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	19	0.17
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	19	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	1	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	1	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	2	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	2	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	4	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	4	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	5	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	5	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	7	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	7	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	11	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	11	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	12	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	12	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	15	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	15	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	17	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	17	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	19	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	19	0.17
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	15	0.17
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	15	0.17
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	13	0.17
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	13	0.17
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	3	0.16
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	3	0.16
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	6	0.16
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	6	0.16
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	8	0.16
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	8	0.16
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	10	0.16
(1,61)	1:B:15:G:H1'	1:B:15:G:H4'	10	0.16
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	8	0.16
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	8	0.16
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	19	0.16
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	19	0.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	2	0.16
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	2	0.16
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	16	0.16
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	16	0.16
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	1	0.16
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	1	0.16
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	11	0.16
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	11	0.16
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	9	0.15
(1,79)	1:B:17:G:H1'	1:B:17:G:H3'	9	0.15
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	5	0.15
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	5	0.15
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	7	0.15
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	7	0.15
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	11	0.15
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	11	0.15
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	2	0.15
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	2	0.15
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	10	0.15
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	10	0.15
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	17	0.15
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	17	0.15
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	19	0.15
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	19	0.15
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	9	0.14
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	9	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	1	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	1	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	4	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	4	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	5	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	5	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	7	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	7	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	8	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	8	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	9	0.14
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	9	0.14
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	1	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	1	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	3	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	3	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	9	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	9	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	10	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	10	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	13	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	13	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	14	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	14	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	18	0.13
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	18	0.13
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	5	0.13
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	5	0.13
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	17	0.13
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	17	0.13
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	18	0.13
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	18	0.13
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	3	0.13
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	3	0.13
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	14	0.13
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	14	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	3	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	3	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	6	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	6	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	14	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	14	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	15	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	15	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	16	0.13
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	16	0.13
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	3	0.12
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	3	0.12
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	9	0.12
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	9	0.12
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	2	0.12
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	2	0.12
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	18	0.12
(1,111)	1:B:21:C:H3'	1:B:21:C:H6	18	0.12
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	6	0.11
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	6	0.11
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	15	0.11
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	15	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	17	0.11
(1,70)	1:B:16:A:H1'	1:B:16:A:H4'	17	0.11
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	15	0.11
(1,56)	1:B:14:C:H3'	1:B:14:C:H6	15	0.11
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	14	0.11
(1,55)	1:B:14:C:H2'	1:B:14:C:H6	14	0.11
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	7	0.11
(1,48)	1:B:13:A:H2'	1:B:14:C:H6	7	0.11

10 Dihedral-angle violation analysis

Dihedral angle analysis failed due to data error in the dihedral angle restraints, possibly missing target value