Full wwPDB NMR Structure Validation Report (i)

Feb 13, 2022 – 11:07 PM EST

PDB ID	:	1HRZ
Title	:	THE 3D STRUCTURE OF THE HUMAN SRY-DNA COMPLEX SOLVED BY MULTI-DIMENSIONAL HETERONUCLEAR-EDITED AND-FILTERED NMR
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Deposited on	:	1995-05-09

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

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

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

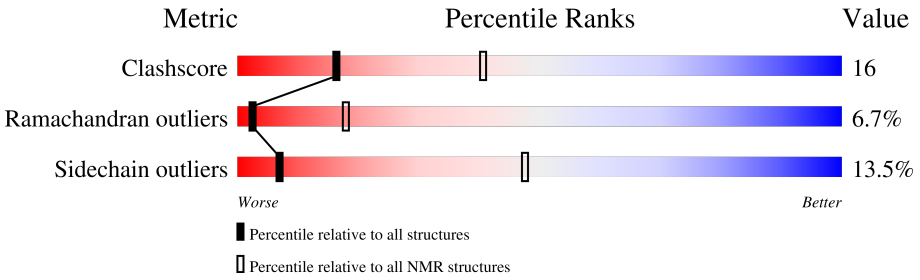
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.26
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.26

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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

Mol	Chain	Length	Quality of chain
1	B	8	38% 62%
2	C	8	12% 50% 38%
3	A	76	50% 38% 5% . .

2 Ensemble composition and analysis

This entry contains 35 models. Model 27 is the overall representative, medoid model (most similar to other models).

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

Well-defined (core) protein residues			
Well-defined core	**Residue range (total)**	**Backbone RMSD (Å)**	**Medoid model**
1	A:5-A:75 (71)	0.42	27

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 9 clusters and 3 single-model clusters were found.

Cluster number	Models
1	18, 22, 23, 27, 28, 29
2	8, 16, 17, 25, 26, 33
3	3, 11, 13, 15, 34
4	6, 14, 30, 35
5	1, 5, 31
6	9, 20
7	12, 24
8	7, 32
9	2, 10
Single-model clusters	4; 19; 21

3 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1794 atoms, of which 828 are hydrogens and 0 are deuteriums.

- Molecule 1 is a DNA chain called DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3').

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
1	B	8	250	77	90	34	42	7	0

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

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
2	C	8	256	79	94	26	50	7	0

- Molecule 3 is a protein called HUMAN SRY.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
3	A	73	1288	409	644	123	107	5	0

4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

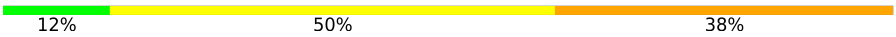
These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for 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: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

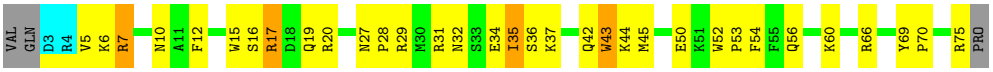
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



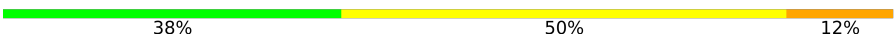
4.2 Scores per residue for each member of the ensemble

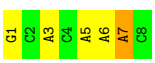
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



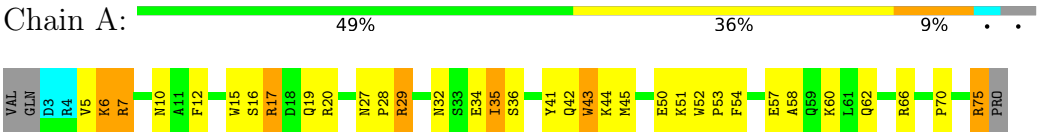


• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

Chain A:



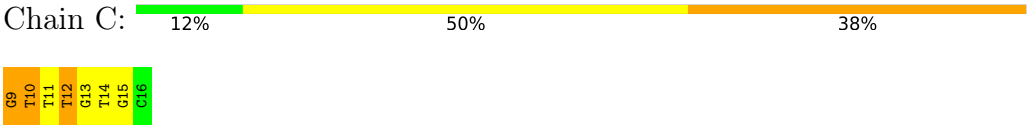
4.2.2 Score per residue for model 2

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

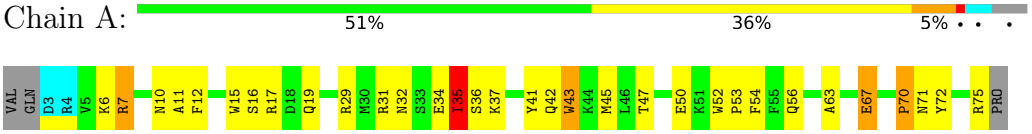
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

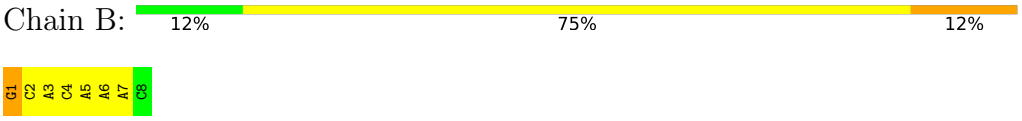
Chain A:



4.2.3 Score per residue for model 3

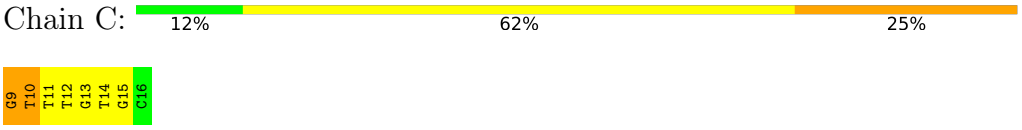
• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



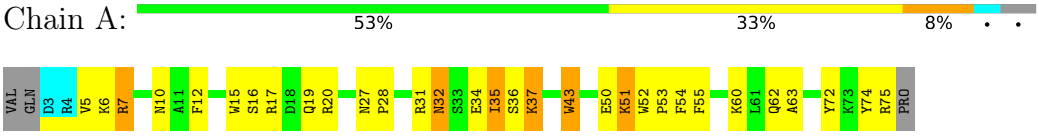
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

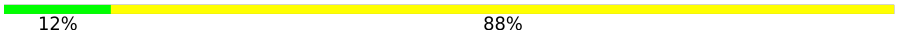
Chain A:



4.2.4 Score per residue for model 4

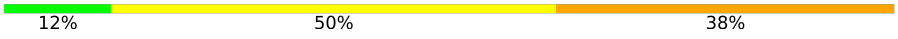
• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



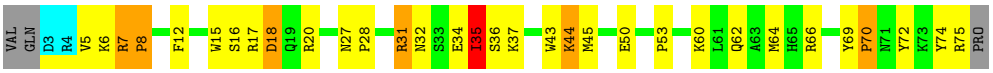
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

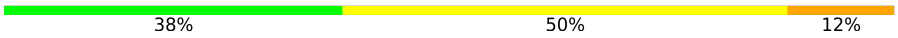
Chain A:



4.2.5 Score per residue for model 5

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

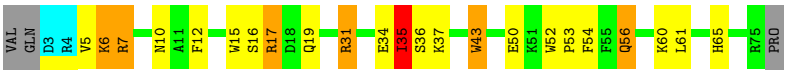


• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

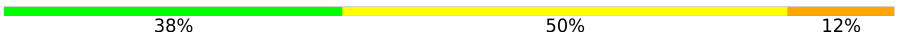
• Molecule 3: HUMAN SRY

Chain A:



4.2.6 Score per residue for model 6

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

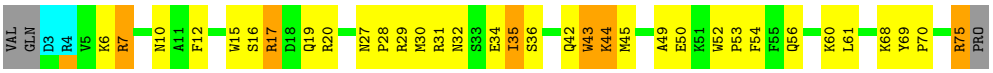
Chain B: 

• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C: 

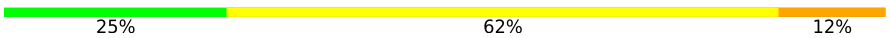
• Molecule 3: HUMAN SRY

Chain A: 



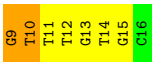
4.2.7 Score per residue for model 7

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B: 

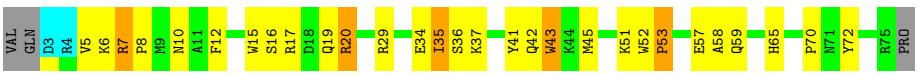
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C: 



• Molecule 3: HUMAN SRY

Chain A: 



4.2.8 Score per residue for model 8

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

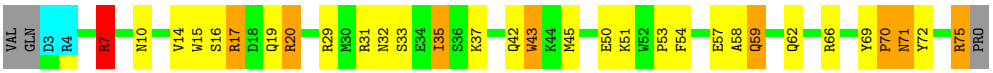
Chain B:

• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

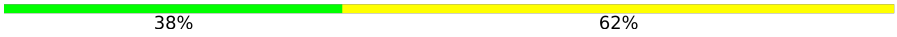
Chain A:



4.2.9 Score per residue for model 9

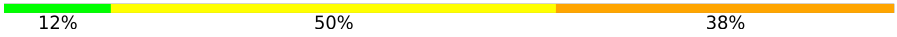
• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



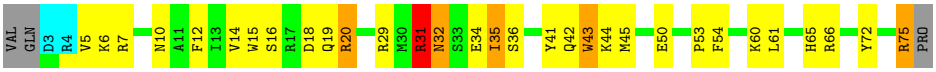
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



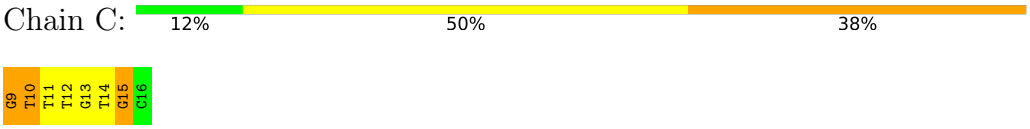
4.2.10 Score per residue for model 10

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

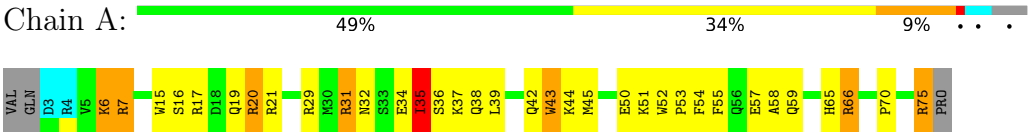
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



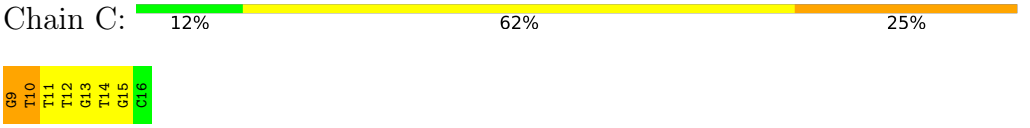
4.2.11 Score per residue for model 11

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

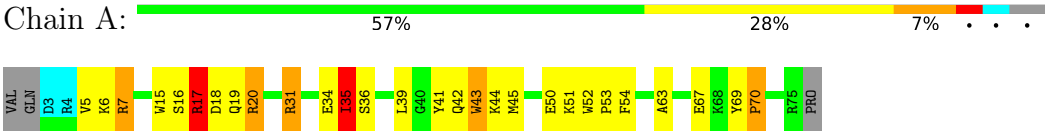
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



4.2.12 Score per residue for model 12

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

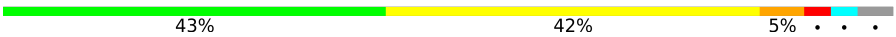
Chain B:

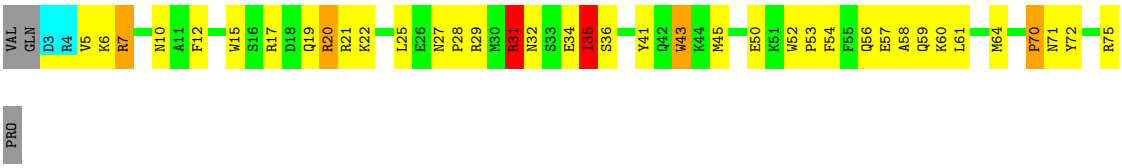
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

Chain A:

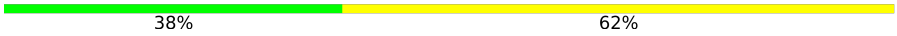




4.2.13 Score per residue for model 13

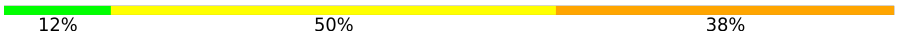
• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



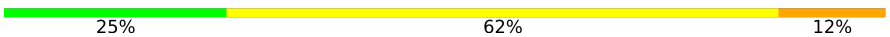
• Molecule 3: HUMAN SRY

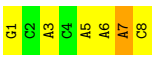
Chain A:

4.2.14 Score per residue for model 14

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

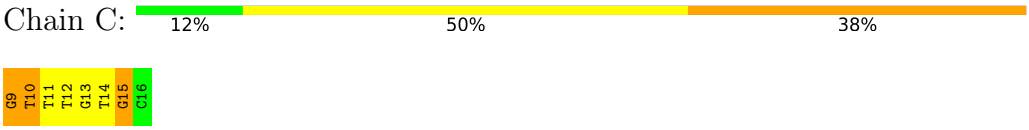
Chain B:





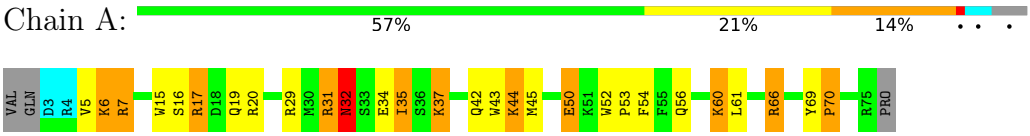
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



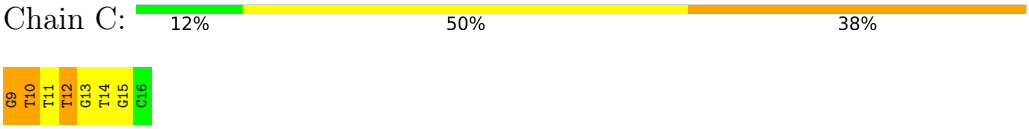
4.2.15 Score per residue for model 15

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

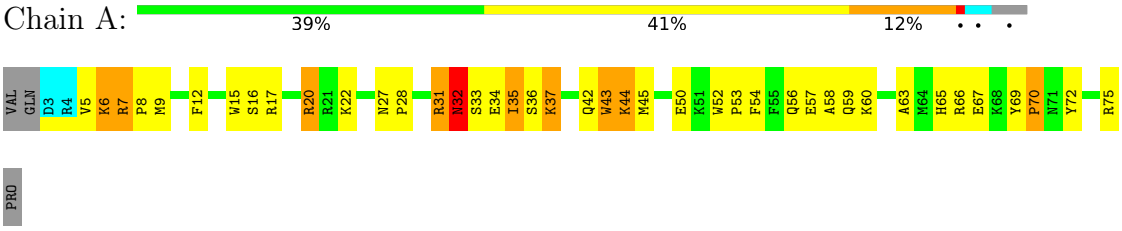
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



4.2.16 Score per residue for model 16

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

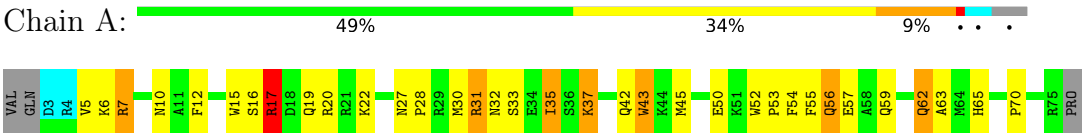
Chain B:

• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

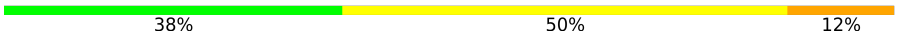
Chain A:



4.2.17 Score per residue for model 17

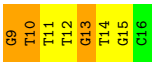
• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



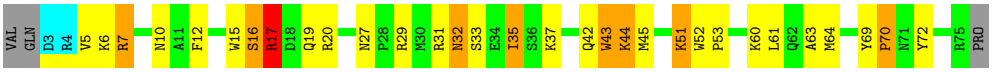
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



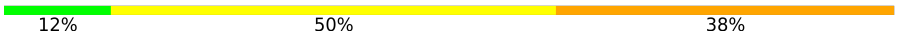
4.2.18 Score per residue for model 18

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

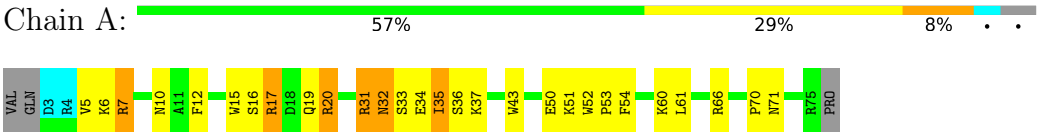
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



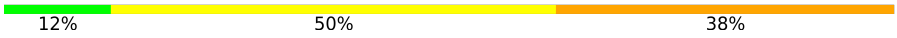
4.2.19 Score per residue for model 19

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

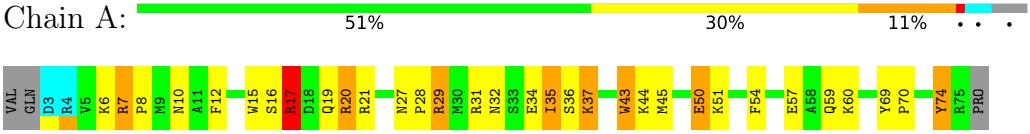
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



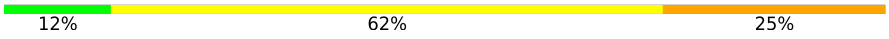
4.2.20 Score per residue for model 20

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

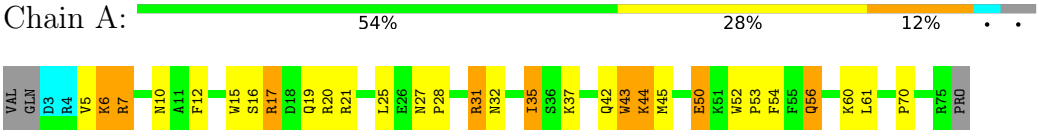
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



4.2.21 Score per residue for model 21

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

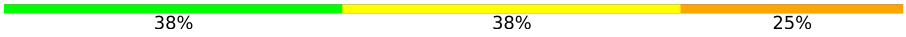
• Molecule 3: HUMAN SRY

Chain A:

4.2.22 Score per residue for model 22

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

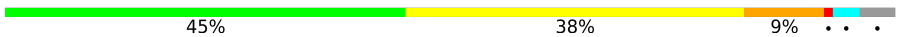


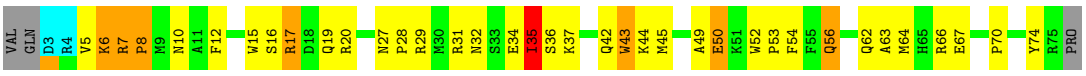
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

Chain A:

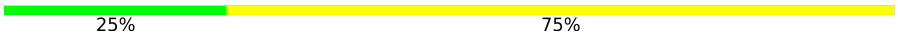


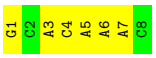


4.2.23 Score per residue for model 23

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



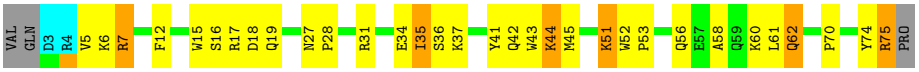


• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

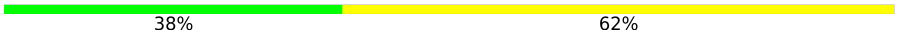
Chain A:



4.2.24 Score per residue for model 24

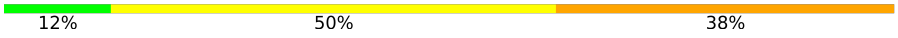
• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



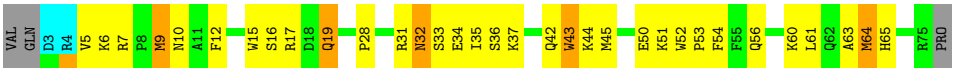
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



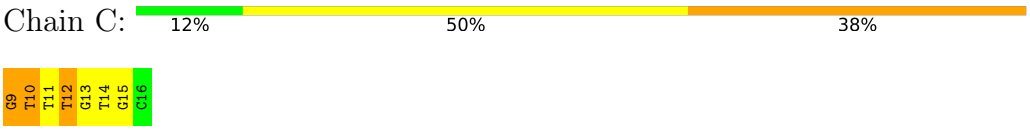
4.2.25 Score per residue for model 25

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

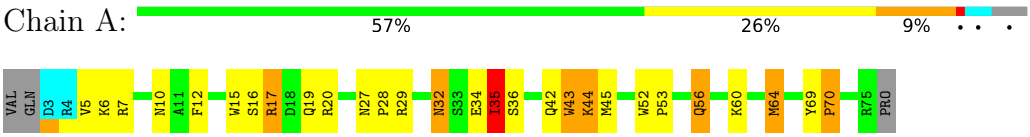
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



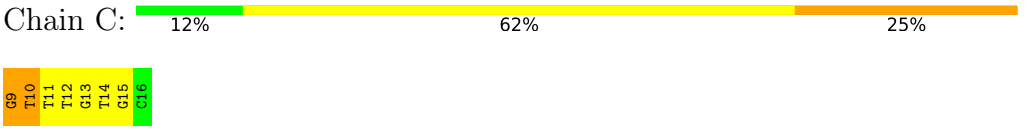
4.2.26 Score per residue for model 26

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

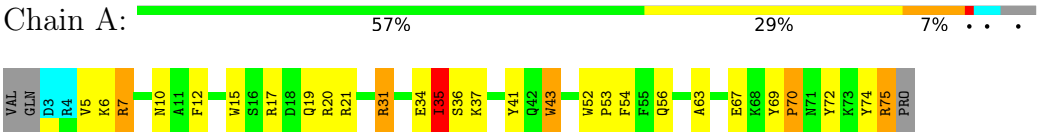
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



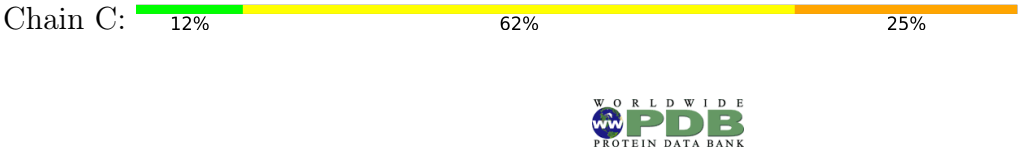
4.2.27 Score per residue for model 27 (medoid)

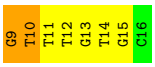
• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

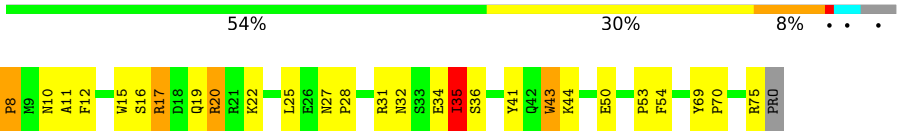
Chain C:





• Molecule 3: HUMAN SRY

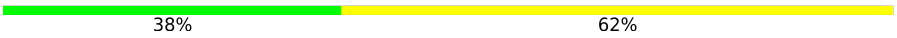
Chain A:



4.2.28 Score per residue for model 28

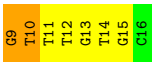
• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



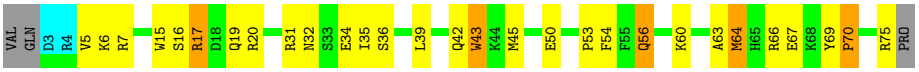
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

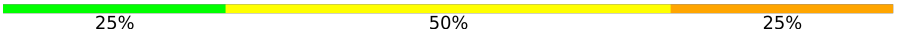
Chain A:

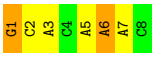


4.2.29 Score per residue for model 29

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:



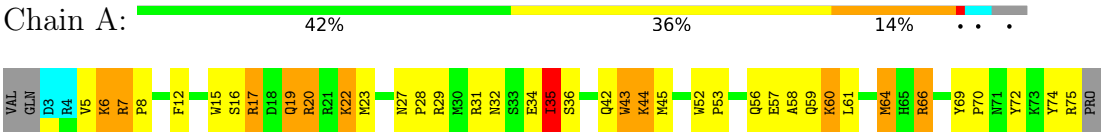


• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

Chain A:



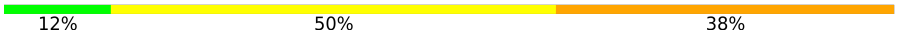
4.2.30 Score per residue for model 30

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

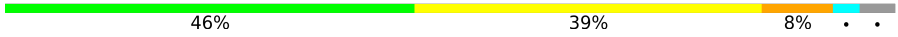
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

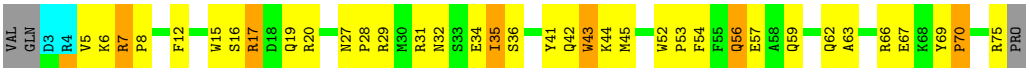
Chain C:



• Molecule 3: HUMAN SRY

Chain A:





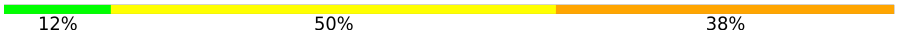
4.2.31 Score per residue for model 31

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

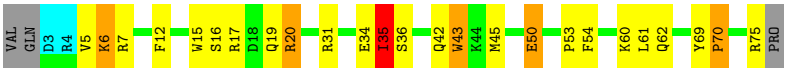
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

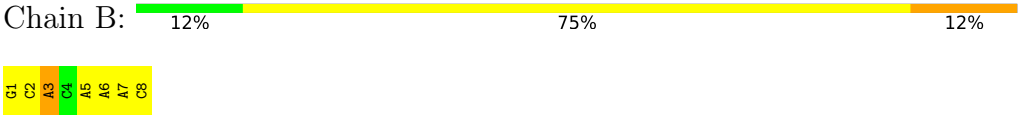
Chain A:



4.2.32 Score per residue for model 32

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

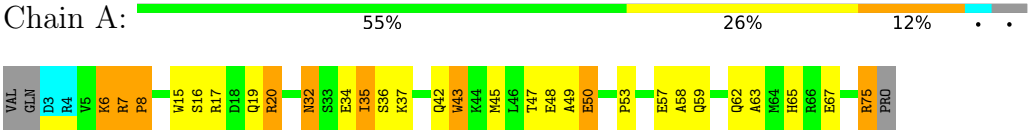


• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

Chain A:



4.2.33 Score per residue for model 33

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

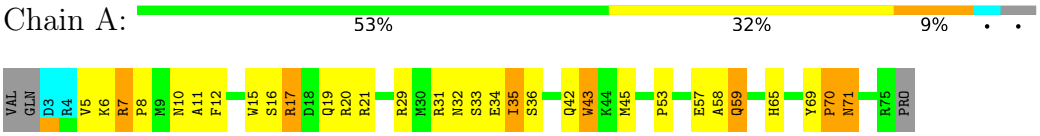
Chain B:

• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:

• Molecule 3: HUMAN SRY

Chain A:



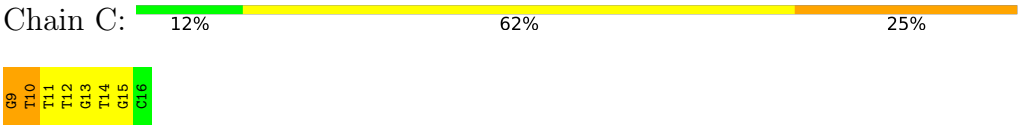
4.2.34 Score per residue for model 34

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

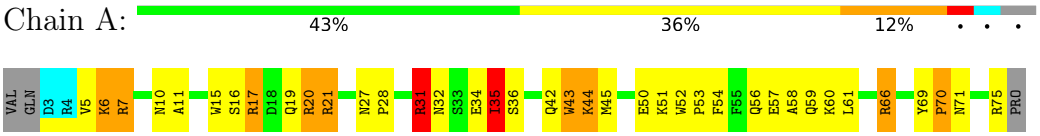
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



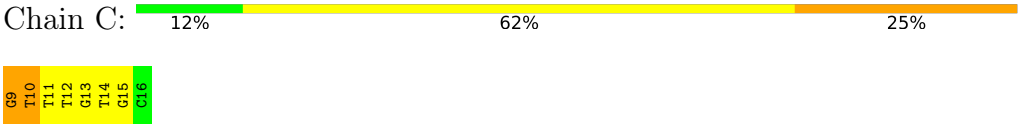
4.2.35 Score per residue for model 35

• Molecule 1: DNA (5'-D(*GP*CP*AP*CP*AP*AP*AP*C)-3')

Chain B:

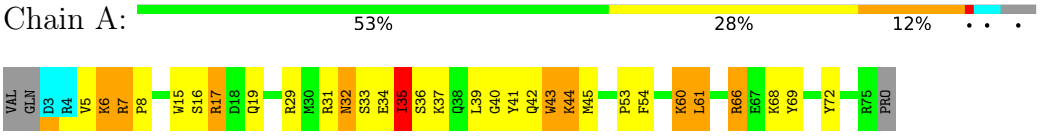
• Molecule 2: DNA (5'-D(*GP*TP*TP*TP*GP*TP*GP*C)-3')

Chain C:



• Molecule 3: HUMAN SRY

Chain A:



5 Refinement protocol and experimental data overview

The models were refined using the following method: *?*.

Of the ? calculated structures, 35 were deposited, based on the following criterion: *?*.

The authors did not provide any information on software used for structure solution, optimization or refinement.

No chemical shift data was provided.

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	B	1.39±0.02	0±0/180 (0.0± 0.1%)	2.88±0.01	16±1/275 (6.0± 0.2%)
2	C	1.55±0.02	3±1/180 (1.8± 0.5%)	2.67±0.01	15±1/277 (5.5± 0.5%)
3	A	1.11±0.01	0±0/642 (0.0± 0.0%)	0.96±0.01	0±0/858 (0.0± 0.0%)
All	All	1.25	118/35070 (0.3%)	1.89	1109/49350 (2.2%)

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	B	1.0±0.0	0.0±0.0
2	C	1.0±0.0	0.0±0.0
3	A	0.0±0.0	0.0±0.2
All	All	70	1

All unique bond 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(Å)	Ideal(Å)	Models	
								Worst	Total
2	C	14	DT	C5-C7	6.52	1.53	1.50	11	35
2	C	11	DT	C5-C7	6.46	1.53	1.50	27	32
2	C	10	DT	C5-C7	6.07	1.53	1.50	8	26
2	C	12	DT	C5-C7	6.00	1.53	1.50	18	23
1	B	1	DG	N9-C8	-5.21	1.34	1.37	8	2

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	1	DG	N7-C8-N9	14.37	120.28	113.10	25	35

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})	Models	
								Worst	Total
2	C	13	DG	N7-C8-N9	14.29	120.24	113.10	3	35
2	C	15	DG	N7-C8-N9	14.17	120.18	113.10	33	35
2	C	9	DG	N7-C8-N9	14.14	120.17	113.10	9	35
1	B	6	DA	N7-C8-N9	12.82	120.21	113.80	7	35
1	B	3	DA	N7-C8-N9	12.73	120.17	113.80	11	35
1	B	7	DA	N7-C8-N9	12.67	120.14	113.80	29	35
1	B	5	DA	N7-C8-N9	12.62	120.11	113.80	25	35
2	C	13	DG	C8-N9-C4	-9.61	102.56	106.40	32	35
2	C	9	DG	C8-N9-C4	-9.49	102.61	106.40	22	35
1	B	1	DG	C8-N9-C4	-9.47	102.61	106.40	21	35
2	C	15	DG	C8-N9-C4	-9.46	102.62	106.40	14	35
1	B	6	DA	C8-N9-C4	-8.90	102.24	105.80	7	35
1	B	7	DA	C8-N9-C4	-8.88	102.25	105.80	14	35
1	B	5	DA	C8-N9-C4	-8.69	102.32	105.80	9	35
1	B	3	DA	C8-N9-C4	-8.47	102.41	105.80	27	35
2	C	9	DG	C5-N7-C8	-8.37	100.12	104.30	7	35
2	C	13	DG	C5-N7-C8	-8.27	100.16	104.30	1	35
1	B	1	DG	C5-N7-C8	-8.27	100.17	104.30	4	35
2	C	15	DG	C5-N7-C8	-8.20	100.20	104.30	16	35
1	B	3	DA	C5-N7-C8	-8.04	99.88	103.90	15	35
1	B	7	DA	C5-N7-C8	-7.80	100.00	103.90	8	35
1	B	5	DA	C5-N7-C8	-7.72	100.04	103.90	4	35
1	B	6	DA	C5-N7-C8	-7.67	100.07	103.90	9	35
1	B	5	DA	O4'-C1'-N9	6.53	112.57	108.00	31	34
2	C	14	DT	O4'-C1'-N1	6.50	112.55	108.00	7	12
2	C	10	DT	C6-C5-C7	-6.38	119.07	122.90	12	34
2	C	12	DT	C6-C5-C7	-6.34	119.10	122.90	5	34
1	B	7	DA	O4'-C1'-N9	6.28	112.39	108.00	17	12
2	C	11	DT	C6-C5-C7	-6.21	119.17	122.90	6	18
2	C	15	DG	O4'-C1'-N9	5.84	112.09	108.00	8	4
1	B	2	DC	O4'-C1'-N1	5.80	112.06	108.00	15	2
2	C	14	DT	C6-C5-C7	-5.72	119.47	122.90	7	4
3	A	7	ARG	NE-CZ-NH1	5.65	123.12	120.30	8	1
2	C	10	DT	C4-C5-C6	5.57	121.34	118.00	11	35
2	C	13	DG	O4'-C1'-N9	5.50	111.85	108.00	34	4
2	C	12	DT	C4-C5-C6	5.43	121.26	118.00	4	33
2	C	12	DT	O4'-C1'-N1	5.39	111.77	108.00	22	5
2	C	9	DG	O4'-C1'-N9	5.36	111.75	108.00	14	12
1	B	3	DA	O4'-C1'-N9	5.33	111.73	108.00	12	1
2	C	11	DT	C4-C5-C6	5.30	121.18	118.00	5	19
2	C	14	DT	C4-C5-C6	5.10	121.06	118.00	33	4

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})	Models	
								Worst	Total
1	B	1	DG	O4'-C1'-N9	5.08	111.56	108.00	9	1

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

Mol	Chain	Res	Type	Atoms	Models (Total)
1	B	5	DA	C3'	35
2	C	9	DG	C3'	35

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
3	A	7	ARG	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
1	B	160	90	90	1±1
2	C	162	94	94	2±1
3	A	625	627	627	26±5
All	All	33145	28385	28385	975

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 16.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:A:19:GLN:NE2	3:A:39:LEU:HD21	0.74	1.97	28	2
3:A:7:ARG:NE	3:A:7:ARG:H	0.70	1.84	14	8
3:A:5:VAL:HG12	3:A:6:LYS:H	0.68	1.46	7	7
3:A:56:GLN:NE2	3:A:56:GLN:N	0.67	2.43	16	9
3:A:56:GLN:H	3:A:56:GLN:NE2	0.66	1.89	28	3
3:A:35:ILE:C	3:A:35:ILE:HD13	0.65	2.12	20	31
3:A:56:GLN:NE2	3:A:56:GLN:H	0.64	1.90	16	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:A:7:ARG:NE	3:A:7:ARG:N	0.64	2.46	3	5
3:A:7:ARG:CG	3:A:7:ARG:HH11	0.63	2.06	8	3
3:A:31:ARG:NE	3:A:31:ARG:H	0.62	1.93	14	1
3:A:15:TRP:CE3	3:A:43:TRP:CH2	0.62	2.87	12	30
3:A:32:ASN:N	3:A:32:ASN:ND2	0.62	2.47	32	2
3:A:7:ARG:HH11	3:A:7:ARG:H	0.61	1.36	15	3
3:A:7:ARG:CD	3:A:7:ARG:N	0.60	2.64	1	11
3:A:64:MET:SD	3:A:64:MET:N	0.60	2.74	29	3
3:A:7:ARG:CG	3:A:7:ARG:NH1	0.60	2.65	8	1
3:A:22:LYS:NZ	3:A:23:MET:N	0.59	2.50	29	1
3:A:5:VAL:HG12	3:A:6:LYS:N	0.59	2.12	5	23
3:A:41:TYR:CE1	3:A:45:MET:SD	0.59	2.95	35	1
3:A:20:ARG:HH11	3:A:20:ARG:CB	0.58	2.12	7	3
3:A:15:TRP:O	3:A:19:GLN:N	0.58	2.37	33	32
3:A:27:ASN:N	3:A:28:PRO:CD	0.58	2.66	23	16
3:A:52:TRP:CD1	3:A:56:GLN:NE2	0.57	2.72	14	3
3:A:31:ARG:NE	3:A:31:ARG:N	0.57	2.51	14	1
3:A:19:GLN:NE2	3:A:22:LYS:NZ	0.57	2.52	12	1
3:A:7:ARG:N	3:A:7:ARG:CD	0.57	2.67	3	4
3:A:22:LYS:HZ2	3:A:23:MET:N	0.57	1.97	29	1
3:A:50:GLU:O	3:A:54:PHE:CE2	0.57	2.58	19	21
3:A:69:TYR:N	3:A:70:PRO:CD	0.57	2.68	14	16
3:A:9:MET:SD	3:A:9:MET:N	0.57	2.77	15	2
3:A:5:VAL:O	3:A:7:ARG:NH1	0.56	2.38	13	7
3:A:20:ARG:HH11	3:A:20:ARG:CG	0.56	2.13	7	3
3:A:34:GLU:OE1	3:A:34:GLU:N	0.56	2.38	22	3
3:A:31:ARG:O	3:A:32:ASN:ND2	0.56	2.39	14	1
2:C:12:DT:O2	3:A:12:PHE:CD2	0.56	2.59	15	19
3:A:50:GLU:O	3:A:54:PHE:CE1	0.56	2.59	18	2
3:A:52:TRP:O	3:A:56:GLN:NE2	0.55	2.38	2	12
3:A:60:LYS:O	3:A:64:MET:SD	0.55	2.65	24	6
3:A:31:ARG:CG	3:A:31:ARG:HH11	0.55	2.14	26	6
3:A:71:ASN:N	3:A:71:ASN:HD22	0.55	2.00	18	4
3:A:5:VAL:O	3:A:7:ARG:CZ	0.55	2.55	11	3
3:A:29:ARG:CG	3:A:29:ARG:HH11	0.54	2.16	8	7
3:A:32:ASN:N	3:A:32:ASN:OD1	0.54	2.41	28	5
3:A:32:ASN:N	3:A:32:ASN:HD22	0.54	2.00	32	2
3:A:14:VAL:HG11	3:A:54:PHE:CD1	0.54	2.37	9	3
3:A:16:SER:OG	3:A:20:ARG:NH2	0.54	2.41	6	2
3:A:17:ARG:CA	3:A:17:ARG:HH11	0.54	2.16	6	1
3:A:50:GLU:O	3:A:54:PHE:CD2	0.54	2.61	19	11

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:A:34:GLU:O	3:A:36:SER:N	0.54	2.40	13	28
1:B:7:DA:C4	1:B:8:DC:C5	0.54	2.96	14	2
3:A:50:GLU:O	3:A:54:PHE:CZ	0.54	2.61	1	2
3:A:50:GLU:N	3:A:50:GLU:OE1	0.54	2.40	3	2
3:A:29:ARG:CG	3:A:29:ARG:NH1	0.54	2.71	8	2
3:A:75:ARG:CG	3:A:75:ARG:NH1	0.53	2.71	1	6
3:A:70:PRO:O	3:A:72:TYR:N	0.53	2.42	12	3
3:A:15:TRP:CD2	3:A:43:TRP:CH2	0.53	2.97	27	9
1:B:1:DG:C5	1:B:2:DC:N4	0.53	2.76	29	4
3:A:7:ARG:HH11	3:A:7:ARG:N	0.53	2.02	34	3
1:B:8:DC:O4'	3:A:20:ARG:NH2	0.53	2.42	4	2
3:A:20:ARG:CG	3:A:20:ARG:NH1	0.53	2.71	13	10
2:C:13:DG:H21	3:A:10:ASN:ND2	0.53	2.02	1	5
3:A:66:ARG:NH1	3:A:66:ARG:CG	0.53	2.70	21	4
3:A:31:ARG:CG	3:A:31:ARG:NH1	0.53	2.71	26	2
3:A:6:LYS:HZ1	3:A:65:HIS:CD2	0.52	2.22	10	1
3:A:10:ASN:N	3:A:10:ASN:ND2	0.52	2.55	13	2
3:A:17:ARG:CG	3:A:17:ARG:HH11	0.52	2.18	17	1
3:A:52:TRP:CD1	3:A:56:GLN:OE1	0.52	2.61	30	1
3:A:31:ARG:NH1	3:A:31:ARG:CG	0.52	2.70	9	5
1:B:8:DC:O3'	3:A:28:PRO:O	0.52	2.27	21	1
3:A:56:GLN:N	3:A:56:GLN:HE21	0.52	2.03	20	3
3:A:8:PRO:CG	3:A:65:HIS:CE1	0.52	2.93	7	1
3:A:29:ARG:NH1	3:A:29:ARG:CG	0.52	2.71	33	6
3:A:75:ARG:CG	3:A:75:ARG:HH11	0.52	2.17	1	4
3:A:66:ARG:CG	3:A:66:ARG:HH11	0.52	2.17	21	4
3:A:17:ARG:CG	3:A:17:ARG:NH1	0.51	2.73	17	1
3:A:58:ALA:O	3:A:62:GLN:NE2	0.51	2.42	23	1
3:A:66:ARG:CG	3:A:66:ARG:NH1	0.51	2.72	34	1
3:A:5:VAL:O	3:A:7:ARG:NE	0.51	2.44	12	1
3:A:20:ARG:CG	3:A:20:ARG:HH11	0.51	2.17	18	5
3:A:52:TRP:NE1	3:A:56:GLN:OE1	0.51	2.43	30	1
3:A:37:LYS:N	3:A:37:LYS:CD	0.51	2.74	14	10
1:B:8:DC:O4'	3:A:20:ARG:NH1	0.51	2.43	32	4
3:A:62:GLN:NE2	3:A:63:ALA:N	0.51	2.59	3	1
3:A:56:GLN:H	3:A:56:GLN:HE21	0.51	1.49	16	1
3:A:65:HIS:CD2	3:A:65:HIS:N	0.51	2.79	5	5
3:A:6:LYS:HZ1	3:A:65:HIS:CG	0.51	2.23	10	1
3:A:41:TYR:CE2	3:A:45:MET:CE	0.51	2.93	1	1
3:A:7:ARG:NH1	3:A:7:ARG:CG	0.50	2.73	7	2
3:A:57:GLU:O	3:A:59:GLN:N	0.50	2.44	7	13

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:A:39:LEU:HD22	3:A:43:TRP:CH2	0.50	2.42	21	3
3:A:52:TRP:CE2	3:A:56:GLN:OE1	0.50	2.64	20	2
3:A:5:VAL:CG1	3:A:6:LYS:N	0.50	2.74	35	15
3:A:31:ARG:O	3:A:32:ASN:O	0.50	2.30	15	4
3:A:18:ASP:N	3:A:18:ASP:OD1	0.50	2.44	23	2
1:B:5:DA:O4'	3:A:7:ARG:NH2	0.50	2.43	8	1
3:A:5:VAL:O	3:A:7:ARG:NH2	0.50	2.45	12	2
3:A:6:LYS:H	3:A:6:LYS:CD	0.50	2.19	32	1
3:A:10:ASN:O	3:A:12:PHE:N	0.50	2.45	19	15
2:C:9:DG:HO5'	2:C:10:DT:P	0.50	2.30	3	6
3:A:47:THR:O	3:A:49:ALA:N	0.50	2.45	32	1
3:A:41:TYR:O	3:A:45:MET:SD	0.50	2.70	12	2
3:A:7:ARG:NH1	3:A:7:ARG:HG2	0.49	2.21	8	2
2:C:9:DG:N2	3:A:33:SER:OG	0.49	2.45	17	1
3:A:17:ARG:N	3:A:17:ARG:HE	0.49	2.04	1	1
3:A:16:SER:OG	3:A:20:ARG:CZ	0.49	2.60	6	2
1:B:6:DA:C4'	3:A:17:ARG:HH21	0.49	2.19	29	1
1:B:7:DA:O4'	3:A:20:ARG:NH1	0.49	2.46	14	2
3:A:31:ARG:N	3:A:31:ARG:CD	0.49	2.75	14	1
3:A:34:GLU:N	3:A:34:GLU:CD	0.49	2.66	31	1
1:B:8:DC:C4'	3:A:20:ARG:HH22	0.49	2.21	26	1
3:A:6:LYS:CD	3:A:6:LYS:N	0.49	2.76	32	1
3:A:27:ASN:N	3:A:28:PRO:HD3	0.49	2.23	12	11
3:A:10:ASN:C	3:A:12:PHE:N	0.49	2.65	25	9
3:A:60:LYS:O	3:A:61:LEU:C	0.49	2.51	17	12
3:A:17:ARG:HH11	3:A:17:ARG:N	0.49	2.06	6	1
1:B:6:DA:C1'	3:A:17:ARG:NH2	0.49	2.76	29	1
1:B:4:DC:C4'	3:A:7:ARG:HH21	0.48	2.21	15	1
3:A:41:TYR:C	3:A:41:TYR:CD1	0.48	2.87	23	3
1:B:7:DA:O4'	3:A:17:ARG:NH2	0.48	2.46	1	1
3:A:75:ARG:NE	3:A:75:ARG:C	0.48	2.67	26	1
3:A:31:ARG:CD	3:A:31:ARG:N	0.48	2.77	20	3
3:A:52:TRP:NE1	3:A:56:GLN:CD	0.48	2.66	26	1
3:A:16:SER:O	3:A:20:ARG:CB	0.48	2.62	20	15
3:A:62:GLN:CD	3:A:63:ALA:N	0.48	2.66	16	2
3:A:44:LYS:C	3:A:45:MET:SD	0.48	2.92	17	19
3:A:75:ARG:HE	3:A:75:ARG:C	0.48	2.10	32	1
3:A:15:TRP:O	3:A:19:GLN:CB	0.48	2.62	23	13
3:A:71:ASN:N	3:A:71:ASN:ND2	0.48	2.60	18	1
3:A:16:SER:O	3:A:17:ARG:C	0.47	2.51	8	25
3:A:17:ARG:N	3:A:17:ARG:NH1	0.47	2.62	6	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:A:35:ILE:C	3:A:35:ILE:CD1	0.47	2.82	8	3
3:A:6:LYS:NZ	3:A:65:HIS:CD2	0.47	2.82	10	1
3:A:65:HIS:N	3:A:65:HIS:CD2	0.47	2.80	24	3
3:A:34:GLU:C	3:A:36:SER:N	0.47	2.68	13	24
3:A:20:ARG:NH1	3:A:20:ARG:CG	0.47	2.78	19	1
3:A:39:LEU:HD22	3:A:43:TRP:CZ2	0.47	2.45	21	1
3:A:7:ARG:NH2	3:A:74:TYR:OH	0.47	2.47	19	2
3:A:62:GLN:O	3:A:66:ARG:CG	0.47	2.61	4	2
3:A:63:ALA:O	3:A:67:GLU:CG	0.47	2.63	26	8
1:B:4:DC:O3'	3:A:7:ARG:NH2	0.47	2.47	15	1
3:A:7:ARG:O	3:A:8:PRO:O	0.46	2.33	4	4
1:B:5:DA:C4'	3:A:7:ARG:HH21	0.46	2.23	8	1
3:A:7:ARG:N	3:A:7:ARG:NE	0.46	2.64	26	3
2:C:13:DG:H21	3:A:10:ASN:HD21	0.46	1.52	17	1
3:A:15:TRP:O	3:A:16:SER:C	0.46	2.54	4	19
3:A:41:TYR:CD1	3:A:41:TYR:C	0.46	2.89	26	6
3:A:51:LYS:O	3:A:52:TRP:C	0.46	2.53	24	9
3:A:32:ASN:ND2	3:A:33:SER:H	0.46	2.09	24	1
1:B:2:DC:N4	1:B:3:DA:N6	0.45	2.64	7	1
3:A:54:PHE:CE1	3:A:55:PHE:CE1	0.45	3.04	16	2
1:B:1:DG:C5	1:B:2:DC:C4	0.45	3.04	29	1
3:A:68:LYS:O	3:A:69:TYR:CD1	0.45	2.70	35	1
2:C:9:DG:C2'	2:C:10:DT:O5'	0.45	2.64	3	35
3:A:42:GLN:O	3:A:45:MET:C	0.45	2.55	15	26
3:A:52:TRP:O	3:A:56:GLN:OE1	0.45	2.35	22	5
3:A:63:ALA:O	3:A:67:GLU:OE1	0.45	2.34	2	1
3:A:70:PRO:C	3:A:72:TYR:H	0.45	2.15	2	2
3:A:15:TRP:CE3	3:A:43:TRP:CZ2	0.45	3.05	21	4
3:A:75:ARG:NE	3:A:75:ARG:H	0.45	2.09	6	1
2:C:13:DG:N2	3:A:10:ASN:OD1	0.45	2.50	5	1
3:A:27:ASN:O	3:A:30:MET:SD	0.45	2.75	6	1
3:A:6:LYS:NZ	3:A:65:HIS:CG	0.45	2.85	10	1
1:B:3:DA:N3	3:A:74:TYR:CD2	0.45	2.85	22	1
3:A:50:GLU:CD	3:A:50:GLU:N	0.44	2.70	32	2
3:A:70:PRO:C	3:A:72:TYR:N	0.44	2.70	2	2
1:B:2:DC:C4	1:B:3:DA:N6	0.44	2.86	32	1
3:A:20:ARG:NH2	3:A:35:ILE:HD12	0.44	2.28	8	1
3:A:73:LYS:O	3:A:74:TYR:O	0.44	2.36	13	1
3:A:32:ASN:CG	3:A:33:SER:N	0.44	2.71	33	3
3:A:20:ARG:CG	3:A:21:ARG:N	0.44	2.81	19	3
1:B:8:DC:C5'	3:A:20:ARG:HH22	0.44	2.25	26	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:A:35:ILE:HD13	3:A:35:ILE:O	0.44	2.12	15	1
3:A:52:TRP:CZ2	3:A:56:GLN:OE1	0.44	2.70	29	2
1:B:1:DG:C6	1:B:2:DC:N4	0.44	2.85	29	1
3:A:62:GLN:NE2	3:A:63:ALA:CA	0.43	2.81	3	1
3:A:49:ALA:HB3	3:A:50:GLU:OE1	0.43	2.13	22	2
3:A:60:LYS:O	3:A:63:ALA:N	0.43	2.51	24	2
1:B:8:DC:C1'	3:A:20:ARG:NH2	0.43	2.81	4	1
3:A:54:PHE:C	3:A:54:PHE:CD1	0.43	2.92	35	3
3:A:31:ARG:O	3:A:32:ASN:OD1	0.43	2.36	3	1
3:A:32:ASN:CG	3:A:33:SER:H	0.43	2.16	16	2
3:A:47:THR:C	3:A:49:ALA:N	0.43	2.72	32	1
3:A:31:ARG:O	3:A:32:ASN:CB	0.43	2.66	14	1
1:B:7:DA:O3'	3:A:20:ARG:NH1	0.43	2.51	31	2
3:A:69:TYR:N	3:A:70:PRO:HD3	0.43	2.29	14	3
3:A:52:TRP:NE1	3:A:56:GLN:NE2	0.43	2.66	24	1
3:A:75:ARG:C	3:A:75:ARG:NE	0.43	2.72	32	1
1:B:8:DC:C1'	3:A:20:ARG:HH22	0.43	2.27	4	1
3:A:32:ASN:OD1	3:A:33:SER:N	0.43	2.52	16	2
3:A:54:PHE:CD1	3:A:54:PHE:C	0.43	2.92	34	2
3:A:7:ARG:O	3:A:7:ARG:CG	0.42	2.66	5	1
3:A:67:GLU:OE1	3:A:67:GLU:N	0.42	2.51	2	1
3:A:7:ARG:H	3:A:7:ARG:CD	0.42	2.27	6	1
3:A:62:GLN:O	3:A:66:ARG:CB	0.42	2.67	22	1
3:A:20:ARG:HH22	3:A:35:ILE:HD12	0.42	1.75	33	1
3:A:57:GLU:C	3:A:59:GLN:N	0.42	2.72	7	4
1:B:7:DA:N3	3:A:20:ARG:NH1	0.42	2.64	22	1
3:A:6:LYS:O	3:A:6:LYS:CG	0.42	2.66	15	1
3:A:19:GLN:HE21	3:A:39:LEU:HD21	0.42	1.66	28	1
1:B:6:DA:C1'	3:A:17:ARG:HH21	0.42	2.27	29	1
3:A:20:ARG:HH11	3:A:20:ARG:HB3	0.42	1.74	10	3
2:C:15:DG:OP1	3:A:66:ARG:NH1	0.42	2.52	10	2
3:A:54:PHE:CE1	3:A:55:PHE:CD2	0.42	3.08	3	1
3:A:52:TRP:CE2	3:A:56:GLN:NE2	0.42	2.88	34	1
3:A:27:ASN:C	3:A:29:ARG:H	0.42	2.19	6	2
2:C:9:DG:O3'	2:C:9:DG:O5'	0.42	2.36	23	2
3:A:60:LYS:C	3:A:64:MET:SD	0.42	2.98	24	1
3:A:32:ASN:O	3:A:34:GLU:N	0.42	2.53	35	1
3:A:32:ASN:HD22	3:A:33:SER:H	0.41	1.58	15	1
3:A:57:GLU:O	3:A:58:ALA:C	0.41	2.58	15	3
3:A:10:ASN:O	3:A:11:ALA:C	0.41	2.57	27	4
3:A:5:VAL:CG2	3:A:72:TYR:CD1	0.41	3.03	9	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:6:DA:C2'	3:A:17:ARG:NH2	0.41	2.83	29	1
3:A:19:GLN:OE1	3:A:19:GLN:C	0.41	2.59	19	2
3:A:5:VAL:CG1	3:A:69:TYR:CZ	0.41	3.04	15	1
2:C:14:DT:C4	2:C:15:DG:O6	0.41	2.73	33	3
3:A:7:ARG:HH11	3:A:7:ARG:HG3	0.41	1.75	8	1
1:B:8:DC:C4'	3:A:20:ARG:HH12	0.41	2.28	30	1
3:A:39:LEU:O	3:A:40:GLY:C	0.41	2.59	35	1
3:A:7:ARG:NH2	3:A:74:TYR:CE1	0.41	2.89	29	1
3:A:43:TRP:N	3:A:43:TRP:CE3	0.41	2.89	21	1
3:A:54:PHE:CZ	3:A:55:PHE:CE1	0.41	3.09	21	1
3:A:62:GLN:N	3:A:62:GLN:OE1	0.41	2.54	23	1
3:A:62:GLN:OE1	3:A:62:GLN:O	0.41	2.39	21	1
3:A:61:LEU:N	3:A:61:LEU:CD2	0.41	2.84	29	1
1:B:7:DA:C1'	3:A:20:ARG:NH1	0.41	2.84	14	1
3:A:7:ARG:H	3:A:7:ARG:HE	0.41	1.57	17	1
3:A:68:LYS:C	3:A:69:TYR:CD1	0.40	2.95	6	1
3:A:10:ASN:N	3:A:10:ASN:HD22	0.40	2.15	13	1
3:A:34:GLU:O	3:A:35:ILE:C	0.40	2.60	3	1
3:A:19:GLN:NE2	3:A:22:LYS:HZ2	0.40	2.13	12	1
3:A:47:THR:OG1	3:A:50:GLU:OE1	0.40	2.39	2	1
3:A:7:ARG:HD3	3:A:7:ARG:N	0.40	2.31	8	1
3:A:20:ARG:HH11	3:A:20:ARG:HG2	0.40	1.76	13	2
3:A:52:TRP:N	3:A:53:PRO:HD2	0.40	2.31	7	1
3:A:5:VAL:HG21	3:A:72:TYR:CE1	0.40	2.51	17	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	
3	A	70/76 (92%)	54±2 (78±3%)	11±2 (16±3%)	5±1 (7±2%)	2	18
All	All	2450/2660 (92%)	1904 (78%)	382 (16%)	164 (7%)	2	18

All 16 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
3	A	53	PRO	34
3	A	17	ARG	25
3	A	70	PRO	25
3	A	35	ILE	16
3	A	32	ASN	13
3	A	8	PRO	11
3	A	31	ARG	10
3	A	51	LYS	7
3	A	58	ALA	7
3	A	74	TYR	5
3	A	71	ASN	3
3	A	25	LEU	3
3	A	61	LEU	2
3	A	16	SER	1
3	A	28	PRO	1
3	A	48	GLU	1

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
3	A	65/70 (93%)	56±2 (86±3%)	9±2 (14±3%)	7 47
All	All	2275/2450 (93%)	1967 (86%)	308 (14%)	7 47

All 30 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)
3	A	7	ARG	35
3	A	35	ILE	35
3	A	43	TRP	35
3	A	31	ARG	21
3	A	6	LYS	20
3	A	20	ARG	18
3	A	75	ARG	18
3	A	37	LYS	17
3	A	44	LYS	13
3	A	66	ARG	12

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Mol	Chain	Res	Type	Models (Total)
3	A	32	ASN	10
3	A	29	ARG	9
3	A	60	LYS	9
3	A	50	GLU	7
3	A	56	GLN	7
3	A	17	ARG	7
3	A	62	GLN	6
3	A	21	ARG	5
3	A	64	MET	5
3	A	22	LYS	4
3	A	18	ASP	3
3	A	59	GLN	2
3	A	38	GLN	2
3	A	19	GLN	2
3	A	67	GLU	1
3	A	10	ASN	1
3	A	30	MET	1
3	A	9	MET	1
3	A	28	PRO	1
3	A	71	ASN	1

6.3.3 RNA

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains

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

6.5 Carbohydrates

There are no monosaccharides in this entry.

6.6 Ligand geometry

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

No chemical shift data were provided