



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 02:01 PM BST

PDB ID : 1AHD
Title : DETERMINATION OF THE NMR SOLUTION STRUCTURE OF AN AN-TENNAPEDIA HOMEODOMAIN-DNA COMPLEX
Authors : Billeter, M.; Qian, Y.Q.; Otting, G.; Muller, M.; Gehring, W.J.; Wuthrich, K.
Deposited on : 1993-04-02

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
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : rb-20027457
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027457

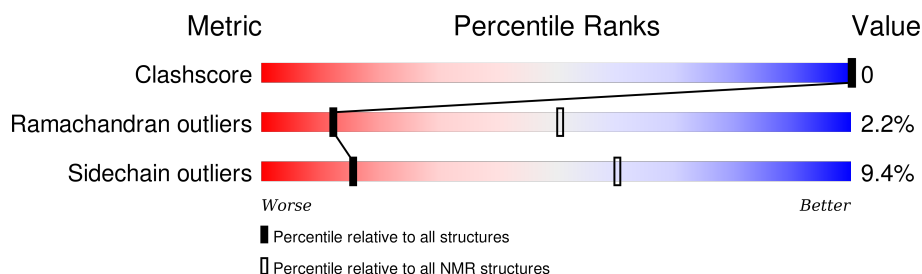
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment 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	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

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	14	
2	B	14	
3	P	68	

2 Ensemble composition and analysis

This entry contains 16 models. Model 16 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	P:7-P:56 (50)	0.48	16

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 2 clusters and 3 single-model clusters were found.

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

3 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2127 atoms, of which 950 are hydrogens and 0 are deuteriums.

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

Mol	Chain	Residues	Atoms						Trace
1	A	14	Total	C	H	N	O	P	0
			447	138	158	60	78	13	

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

Mol	Chain	Residues	Atoms						Trace
2	B	14	Total	C	H	N	O	P	0
			441	136	162	44	86	13	

- Molecule 3 is a protein called Homeotic protein antennapedia.

Mol	Chain	Residues	Atoms						Trace
3	P	68	Total	C	H	N	O	S	0
			1239	383	630	125	99	2	

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	0	MET	-	EXPRESSION TAG	UNP Q24645
P	39	SER	CYS	ENGINEERED MUTATION	UNP Q24645

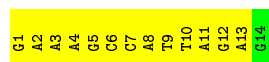
4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA 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*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')

Chain A: 



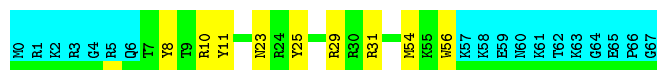
- Molecule 2: DNA (5'-D(*CP*TP*CP*TP*AP*AP*TP*GP*GP*CP*TP*TP*TP*C)-3')

Chain B: 



- Molecule 3: Homeotic protein antennapedia

Chain P: 



4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

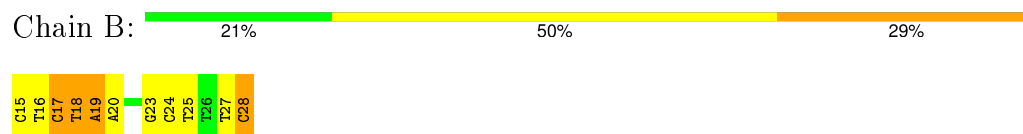
4.2.1 Score per residue for model 1

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')

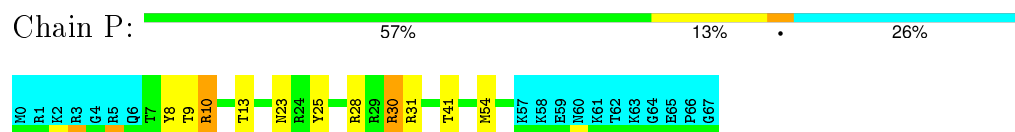
Chain A: 



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

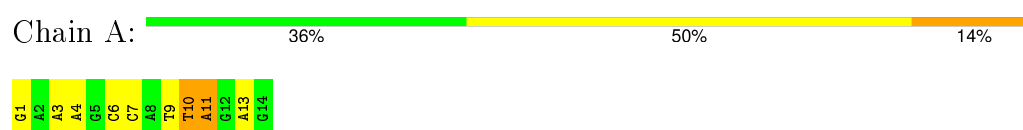


- Molecule 3: Homeotic protein antennapedia



4.2.2 Score per residue for model 2

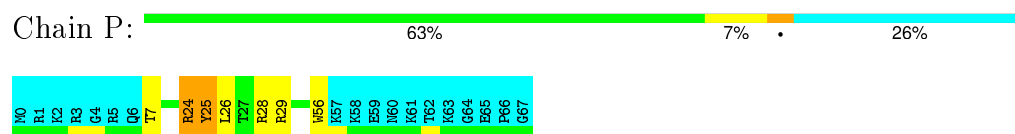
- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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



- Molecule 3: Homeotic protein antennapedia



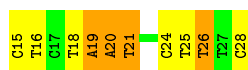
4.2.3 Score per residue for model 3

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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



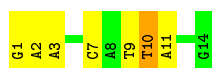


- Molecule 3: Homeotic protein antennapedia

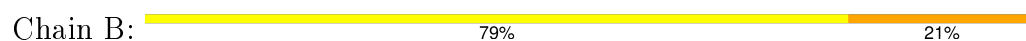


4.2.4 Score per residue for model 4

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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



- Molecule 3: Homeotic protein antennapedia

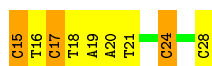


4.2.5 Score per residue for model 5

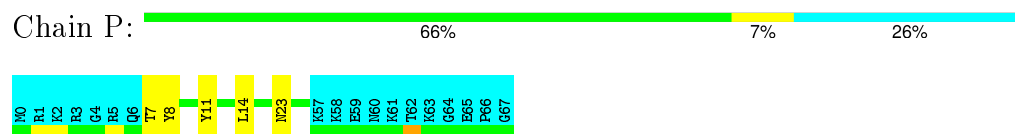
- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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



- Molecule 3: Homeotic protein antennapedia



4.2.6 Score per residue for model 6

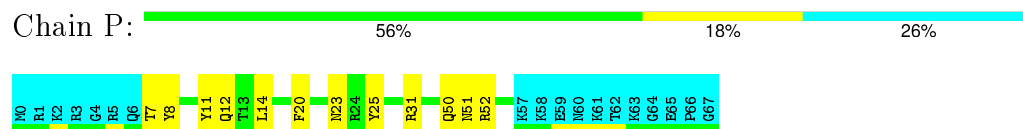
- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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



- Molecule 3: Homeotic protein antennapedia

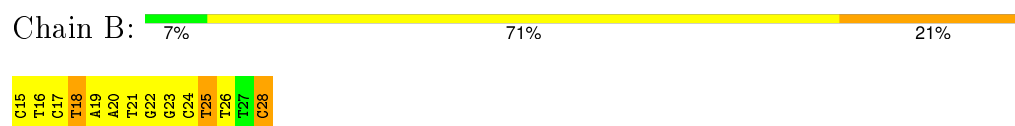


4.2.7 Score per residue for model 7

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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



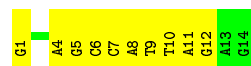
- Molecule 3: Homeotic protein antennapedia



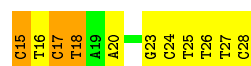


4.2.8 Score per residue for model 8

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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



- Molecule 3: Homeotic protein antennapedia

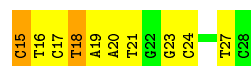


4.2.9 Score per residue for model 9

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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



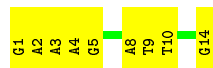
- Molecule 3: Homeotic protein antennapedia



4.2.10 Score per residue for model 10

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')

Chain A: 



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

Chain B: 



- Molecule 3: Homeotic protein antennapedia

Chain P: 



4.2.11 Score per residue for model 11

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')

Chain A: 



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

Chain B: 



- Molecule 3: Homeotic protein antennapedia

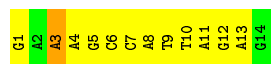
Chain P: 



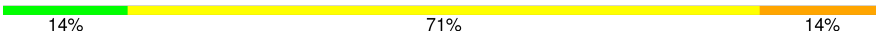
4.2.12 Score per residue for model 12

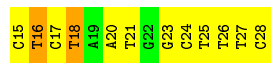
- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')

Chain A: 



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

Chain B: 



- Molecule 3: Homeotic protein antennapedia

Chain P: 



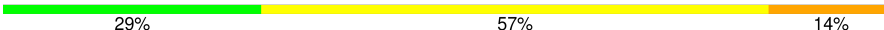
4.2.13 Score per residue for model 13

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')

Chain A: 



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

Chain B: 



- Molecule 3: Homeotic protein antennapedia

Chain P: 



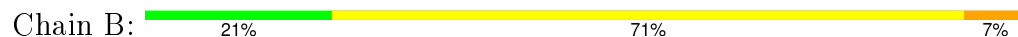
4.2.14 Score per residue for model 14

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')

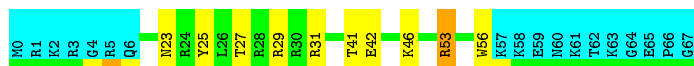
Chain A: 



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

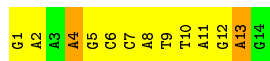


- Molecule 3: Homeotic protein antennapedia

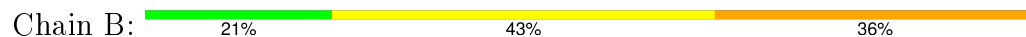


4.2.15 Score per residue for model 15

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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

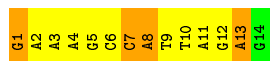
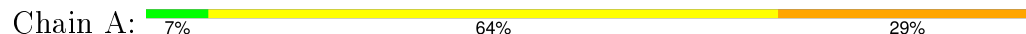


- Molecule 3: Homeotic protein antennapedia



4.2.16 Score per residue for model 16 (medoid)

- Molecule 1: DNA (5'-D(*GP*AP*AP*AP*GP*CP*CP*AP*TP*TP*AP*GP*AP*G)-3')



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

Chain B:  14% 79% 7%

C15
T16
C17
T18
A19
A20
T21
G22
G23
C24
T25
T26
T27
C28

- Molecule 3: Homeotic protein antennapedia

Chain P:  63% 7% 26%

H0
R1
R2
R3
G4
R5
G6
T7
Y8
F22
Y25
T41
R52
R53
R54
R57
R58
R59
H60
R61
R62
R63
G64
R65
P66
G67

5 Refinement protocol and experimental data overview ⓘ

Of the ? calculated structures, 16 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. No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality

6.1 Standard geometry

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	1.26±0.02	0±0/326 (0.0±0.0%)	1.96±0.07	10±3/502 (1.9±0.6%)
2	B	1.30±0.02	0±0/310 (0.0±0.0%)	2.12±0.06	13±3/476 (2.7±0.6%)
3	P	0.63±0.01	0±0/474 (0.0±0.0%)	1.05±0.05	1±1/636 (0.2±0.1%)
All	All	1.05	0/17760 (0.0%)	1.72	377/25824 (1.5%)

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	4.4±1.5
2	B	0.0±0.0	5.3±1.9
3	P	0.0±0.0	3.8±1.5
All	All	0	217

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
2	B	15	DC	O4'-C1'-N1	11.93	116.35	108.00	5	16
2	B	18	DT	O4'-C1'-N1	11.68	116.17	108.00	5	4
1	A	11	DA	O4'-C1'-N9	11.29	115.90	108.00	7	9
2	B	16	DT	O4'-C1'-N1	10.74	115.52	108.00	5	9
1	A	1	DG	O4'-C1'-N9	9.88	114.91	108.00	5	11
2	B	20	DA	O4'-C1'-N9	9.79	114.86	108.00	12	10
2	B	21	DT	O4'-C1'-N1	8.83	114.18	108.00	9	3
1	A	4	DA	O4'-C1'-N9	8.35	113.85	108.00	13	6
1	A	6	DC	O4'-C1'-N1	8.19	113.73	108.00	3	4
1	A	7	DC	O4'-C1'-N1	8.16	113.72	108.00	5	7
1	A	4	DA	O4'-C4'-C3'	-7.79	101.32	106.00	15	4

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	26	DT	O4'-C1'-N1	7.70	113.39	108.00	14	8
1	A	1	DG	O4'-C1'-C2'	-7.54	99.87	105.90	5	1
2	B	18	DT	O4'-C1'-C2'	-7.23	100.11	105.90	5	1
2	B	27	DT	C6-C5-C7	-7.23	118.56	122.90	1	10
2	B	19	DA	O4'-C1'-N9	7.17	113.02	108.00	15	2
2	B	28	DC	O4'-C1'-N1	7.02	112.92	108.00	2	2
1	A	13	DA	O4'-C4'-C3'	7.02	110.21	106.00	6	4
2	B	20	DA	O4'-C1'-C2'	-6.90	100.38	105.90	13	9
2	B	18	DT	C6-C5-C7	-6.87	118.78	122.90	8	5
2	B	25	DT	O4'-C1'-N1	6.84	112.79	108.00	7	3
2	B	18	DT	C1'-O4'-C4'	-6.84	103.26	110.10	1	5
3	P	56	TRP	C-N-CA	6.83	138.79	121.70	12	2
1	A	5	DG	O4'-C1'-N9	6.81	112.77	108.00	5	2
1	A	12	DG	O4'-C1'-N9	6.80	112.76	108.00	7	4
1	A	7	DC	N1-C2-O2	6.80	122.98	118.90	16	13
2	B	17	DC	O4'-C1'-N1	6.73	112.71	108.00	5	3
3	P	52	ARG	NE-CZ-NH2	-6.68	116.96	120.30	13	1
1	A	9	DT	C6-C5-C7	-6.59	118.95	122.90	14	14
2	B	25	DT	C6-C5-C7	-6.59	118.95	122.90	10	9
2	B	25	DT	O4'-C1'-C2'	-6.58	100.63	105.90	7	2
1	A	10	DT	C6-C5-C7	-6.54	118.98	122.90	6	12
2	B	17	DC	O4'-C4'-C3'	6.53	109.92	106.00	10	1
2	B	24	DC	N1-C2-O2	6.48	122.79	118.90	1	11
2	B	16	DT	C6-C5-C7	-6.39	119.07	122.90	4	8
3	P	25	TYR	CB-CG-CD1	-6.38	117.17	121.00	16	1
1	A	5	DG	O4'-C1'-C2'	-6.33	100.84	105.90	5	1
1	A	1	DG	C4'-C3'-C2'	-6.31	97.42	103.10	2	1
1	A	3	DA	O4'-C1'-N9	6.29	112.41	108.00	12	3
1	A	5	DG	C4'-C3'-C2'	-6.27	97.46	103.10	11	1
1	A	4	DA	O4'-C1'-C2'	-6.26	100.89	105.90	3	2
2	B	15	DC	N1-C2-O2	6.26	122.65	118.90	14	6
2	B	21	DT	O4'-C4'-C3'	-6.25	102.00	104.50	13	1
1	A	12	DG	C1'-O4'-C4'	-6.19	103.91	110.10	1	2
1	A	3	DA	C4'-C3'-C2'	-6.18	97.54	103.10	12	1
1	A	12	DG	P-O3'-C3'	6.18	127.11	119.70	3	1
2	B	15	DC	C1'-O4'-C4'	-6.13	103.97	110.10	11	3
2	B	28	DC	N3-C2-O2	-6.12	117.62	121.90	7	8
3	P	53	ARG	CD-NE-CZ	6.12	132.17	123.60	14	1
1	A	2	DA	O4'-C1'-N9	6.10	112.27	108.00	16	2
2	B	26	DT	C6-C5-C7	-6.10	119.24	122.90	8	10
1	A	6	DC	N1-C2-O2	6.05	122.53	118.90	5	10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	9	DT	O4'-C1'-N1	6.02	112.21	108.00	7	8
2	B	15	DC	N3-C2-O2	-6.00	117.70	121.90	14	4
2	B	20	DA	P-O3'-C3'	5.98	126.88	119.70	5	4
2	B	26	DT	C1'-O4'-C4'	-5.97	104.13	110.10	14	2
2	B	24	DC	N3-C2-O2	-5.95	117.73	121.90	1	4
1	A	7	DC	C1'-O4'-C4'	-5.95	104.15	110.10	8	1
2	B	19	DA	C4-C5-C6	-5.91	114.05	117.00	3	1
3	P	31	ARG	NE-CZ-NH2	-5.88	117.36	120.30	13	1
3	P	56	TRP	CA-CB-CG	5.87	124.86	113.70	7	1
2	B	15	DC	O4'-C1'-C2'	-5.86	101.21	105.90	10	1
3	P	28	ARG	NE-CZ-NH2	-5.84	117.38	120.30	7	2
2	B	24	DC	C1'-O4'-C4'	-5.77	104.33	110.10	4	1
2	B	17	DC	N1-C2-O2	5.76	122.36	118.90	15	5
2	B	19	DA	P-O3'-C3'	5.75	126.61	119.70	10	5
1	A	8	DA	O4'-C4'-C3'	5.74	109.45	106.00	6	5
1	A	11	DA	P-O3'-C3'	5.74	126.58	119.70	6	1
2	B	16	DT	C1'-O4'-C4'	-5.72	104.38	110.10	5	1
3	P	29	ARG	NE-CZ-NH2	-5.71	117.44	120.30	2	1
1	A	1	DG	C1'-O4'-C4'	-5.70	104.40	110.10	5	2
1	A	8	DA	C5-C6-N6	-5.70	119.14	123.70	1	3
2	B	28	DC	N1-C2-O2	5.69	122.31	118.90	5	8
2	B	17	DC	N3-C2-O2	-5.62	117.97	121.90	4	2
2	B	20	DA	C1'-O4'-C4'	-5.58	104.52	110.10	10	3
2	B	17	DC	P-O3'-C3'	5.58	126.39	119.70	16	2
1	A	3	DA	C1'-O4'-C4'	-5.49	104.61	110.10	14	1
1	A	2	DA	O4'-C4'-C3'	5.47	109.28	106.00	15	2
2	B	21	DT	C6-C5-C7	-5.46	119.63	122.90	3	3
3	P	25	TYR	CB-CG-CD2	-5.41	117.75	121.00	8	1
1	A	6	DC	O4'-C4'-C3'	5.37	109.22	106.00	8	1
3	P	24	ARG	NE-CZ-NH2	-5.36	117.62	120.30	2	2
2	B	27	DT	N3-C2-O2	-5.35	119.09	122.30	10	1
1	A	7	DC	N3-C2-O2	-5.34	118.16	121.90	9	4
2	B	16	DT	O4'-C1'-C2'	-5.30	101.66	105.90	5	1
1	A	9	DT	O4'-C1'-C2'	-5.29	101.66	105.90	2	2
2	B	15	DC	C4'-C3'-C2'	-5.26	98.36	103.10	3	1
3	P	53	ARG	NE-CZ-NH2	-5.26	117.67	120.30	9	1
3	P	8	TYR	CB-CG-CD2	-5.26	117.84	121.00	11	2
2	B	27	DT	O4'-C1'-N1	5.23	111.66	108.00	1	1
1	A	8	DA	O4'-C1'-N9	5.21	111.65	108.00	16	1
2	B	19	DA	O4'-C1'-C2'	-5.21	101.73	105.90	7	1
2	B	24	DC	O4'-C1'-C2'	-5.20	101.74	105.90	8	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	14	DG	C4'-C3'-C2'	-5.20	98.42	103.10	13	1
2	B	28	DC	C1'-O4'-C4'	-5.16	104.94	110.10	8	2
1	A	11	DA	C1'-O4'-C4'	-5.16	104.94	110.10	7	1
2	B	25	DT	C1'-O4'-C4'	-5.13	104.97	110.10	8	1
1	A	11	DA	O4'-C1'-C2'	-5.09	101.83	105.90	7	1
2	B	25	DT	N3-C2-O2	-5.07	119.26	122.30	8	1
2	B	20	DA	O4'-C4'-C3'	5.07	109.04	106.00	16	1
1	A	13	DA	P-O3'-C3'	5.06	125.77	119.70	2	2
3	P	10	ARG	NE-CZ-NH2	-5.05	117.77	120.30	15	1
1	A	8	DA	O4'-C1'-C2'	-5.05	101.86	105.90	10	1
2	B	18	DT	C4'-C3'-C2'	-5.03	98.57	103.10	11	1
2	B	27	DT	O4'-C1'-C2'	-5.01	101.89	105.90	8	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)
2	B	18	DT	Sidechain	15
2	B	23	DG	Sidechain	11
3	P	25	TYR	Sidechain	11
1	A	3	DA	Sidechain	10
2	B	24	DC	Sidechain	9
3	P	8	TYR	Sidechain	9
1	A	5	DG	Sidechain	8
3	P	11	TYR	Sidechain	8
2	B	21	DT	Sidechain	8
1	A	12	DG	Sidechain	8
1	A	10	DT	Sidechain	8
2	B	17	DC	Sidechain	7
1	A	4	DA	Sidechain	7
3	P	29	ARG	Sidechain,Peptide	6
2	B	22	DG	Sidechain	5
2	B	28	DC	Sidechain	5
1	A	2	DA	Sidechain	5
2	B	27	DT	Sidechain	5
1	A	13	DA	Sidechain	5
3	P	31	ARG	Sidechain	5
1	A	11	DA	Sidechain	4
2	B	20	DA	Sidechain	4
2	B	19	DA	Sidechain	4

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Mol	Chain	Res	Type	Group	Models (Total)
1	A	1	DG	Sidechain	4
2	B	15	DC	Sidechain	4
2	B	25	DT	Sidechain	4
3	P	10	ARG	Sidechain	4
1	A	6	DC	Sidechain	3
3	P	20	PHE	Sidechain	3
3	P	53	ARG	Sidechain	3
3	P	30	ARG	Sidechain	3
1	A	9	DT	Sidechain	3
3	P	28	ARG	Sidechain	3
2	B	16	DT	Sidechain	2
2	B	26	DT	Sidechain	2
1	A	7	DC	Sidechain	2
3	P	22	PHE	Sidechain	2
1	A	8	DA	Sidechain	2
1	A	14	DG	Sidechain	2
3	P	43	ARG	Sidechain	1
3	P	24	ARG	Sidechain	1
3	P	52	ARG	Sidechain	1

6.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	289	158	158	0±0
2	B	279	162	162	0±0
All	All	16480	12608	12608	1

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 0.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:11:DA:C2	2:B:19:DA:C2	0.42	3.08	1	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	P	50/68 (74%)	43±2 (85±4%)	6±2 (13±3%)	1±1 (2±2%)	12	51
All	All	800/1088 (74%)	681 (85%)	101 (13%)	18 (2%)	12	51

All 13 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
3	P	10	ARG	3
3	P	9	THR	3
3	P	28	ARG	2
3	P	27	THR	1
3	P	56	TRP	1
3	P	24	ARG	1
3	P	29	ARG	1
3	P	21	HIS	1
3	P	7	THR	1
3	P	54	MET	1
3	P	26	LEU	1
3	P	30	ARG	1
3	P	40	LEU	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	P	48/63 (76%)	44±2 (91±4%)	5±2 (9±4%)	15	60
All	All	768/1008 (76%)	696 (91%)	72 (9%)	15	60

All 25 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	P	23	ASN	12
3	P	54	MET	7
3	P	14	LEU	5
3	P	56	TRP	5
3	P	7	THR	4
3	P	13	THR	4
3	P	9	THR	4
3	P	42	GLU	4
3	P	52	ARG	3
3	P	41	THR	3
3	P	31	ARG	3
3	P	24	ARG	2
3	P	25	TYR	2
3	P	39	SER	2
3	P	16	LEU	2
3	P	50	GLN	1
3	P	51	ASN	1
3	P	10	ARG	1
3	P	26	LEU	1
3	P	30	ARG	1
3	P	46	LYS	1
3	P	53	ARG	1
3	P	27	THR	1
3	P	22	PHE	1
3	P	12	GLN	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 carbohydrates in this entry.

6.6 Ligand geometry

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided