



Full wwPDB NMR Structure Validation Report ⓘ

Mar 5, 2026 – 09:14 PM UTC

PDB ID : 7ZEW / pdb_00007zew
BMRB ID : 34725
Title : Complex Cyp33-RRM : AAUAAA RNA
Authors : Blatter, M.; Allain, F.; Meylan, C.
Deposited on : 2022-03-31

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 : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

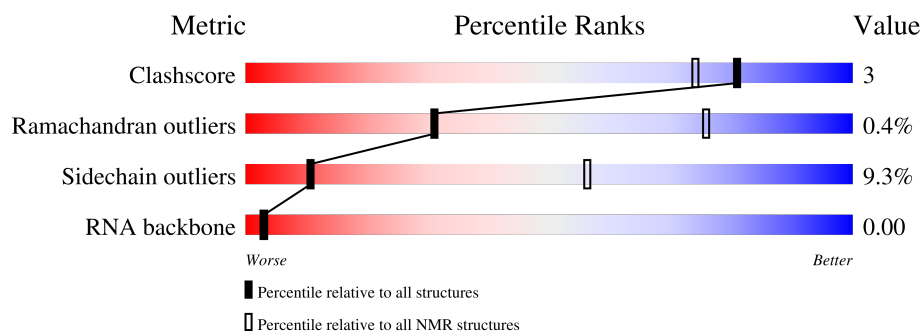
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 78%.

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



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823
RNA backbone	8273	777

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	117	
2	B	6	

2 Ensemble composition and analysis

This entry contains 20 models. Model 1 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:83, A:96-A:105 (89)	0.32	1

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 1 single-model cluster was found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2028 atoms, of which 973 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Peptidyl-prolyl cis-trans isomerase E.

Mol	Chain	Residues	Atoms						Trace
1	A	117	Total	C	H	N	O	S	0
			1836	589	908	157	179	3	

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ALA	-	expression tag	UNP Q9UNP9
A	-2	GLY	-	expression tag	UNP Q9UNP9
A	-1	HIS	-	expression tag	UNP Q9UNP9

- Molecule 2 is a RNA chain called RNA (5'-R(*AP*AP*UP*AP*AP*A)-3').

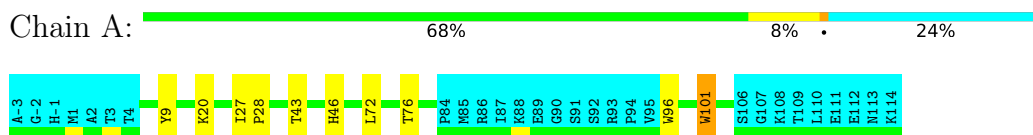
Mol	Chain	Residues	Atoms						Trace
2	B	6	Total	C	H	N	O	P	0
			192	59	65	27	36	5	

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: Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')

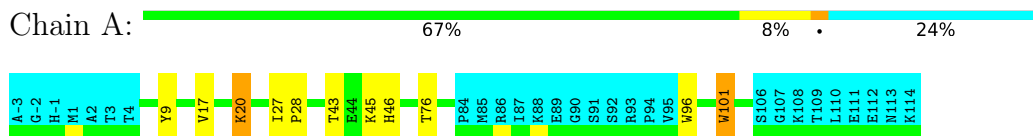


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1 (medoid)

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

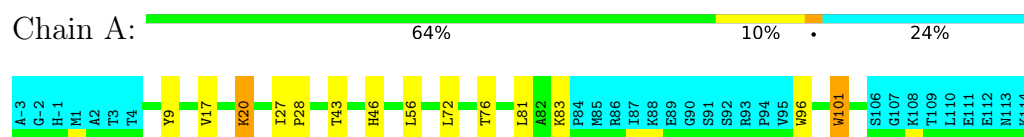


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.2 Score per residue for model 2

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

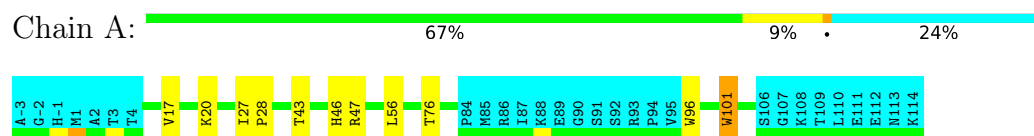


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.3 Score per residue for model 3

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

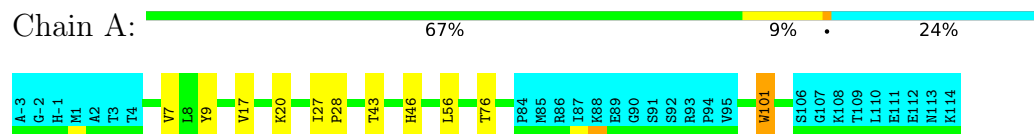


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.4 Score per residue for model 4

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

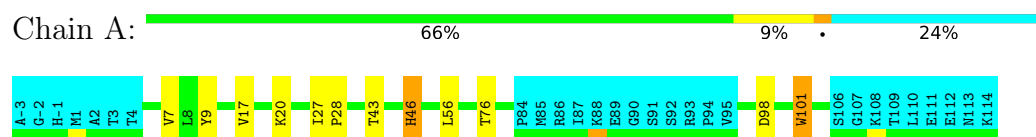


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.5 Score per residue for model 5

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

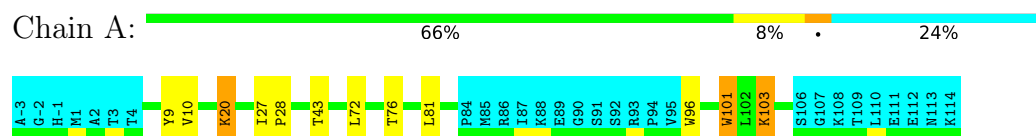


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.6 Score per residue for model 6

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

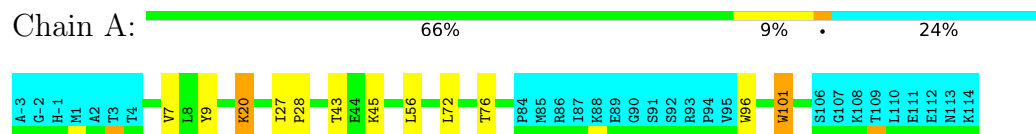


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.7 Score per residue for model 7

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

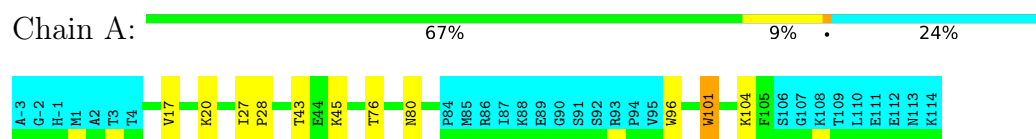


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.8 Score per residue for model 8

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

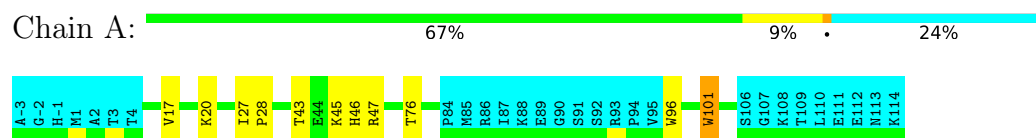


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.9 Score per residue for model 9

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

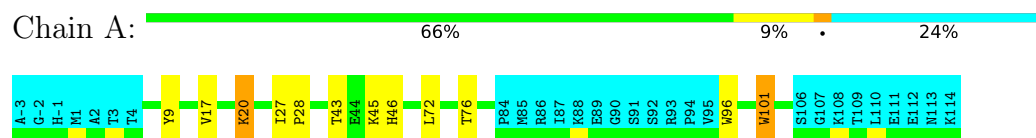


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.10 Score per residue for model 10

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

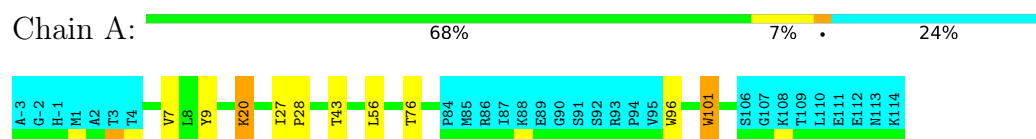


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.11 Score per residue for model 11

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

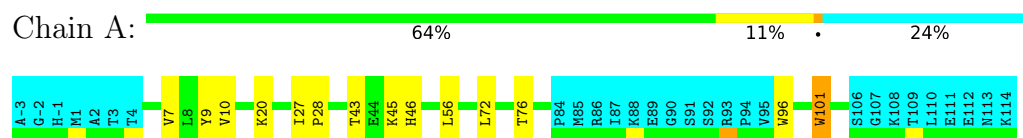


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.12 Score per residue for model 12

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

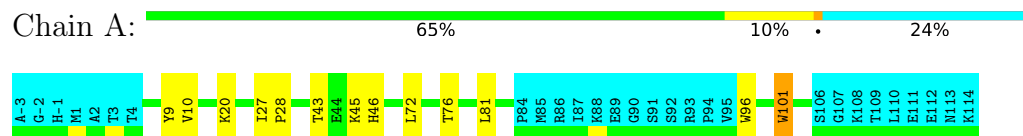


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.13 Score per residue for model 13

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

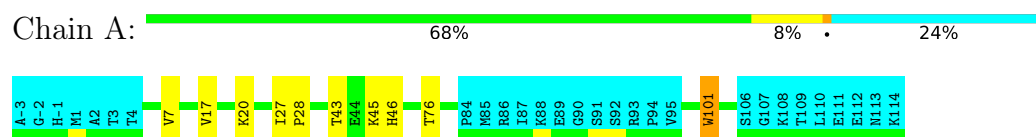


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.14 Score per residue for model 14

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

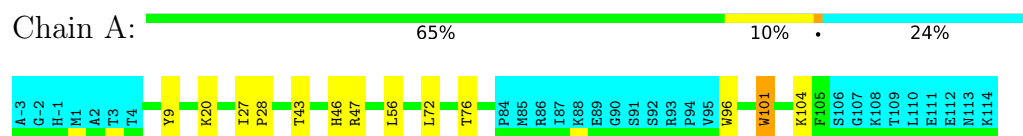


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.15 Score per residue for model 15

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

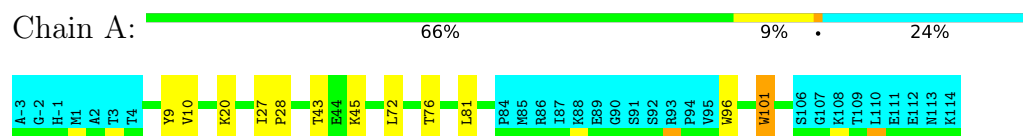


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.16 Score per residue for model 16

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

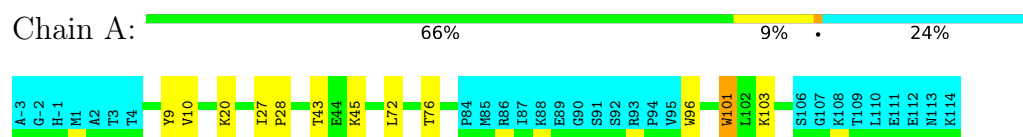


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.17 Score per residue for model 17

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

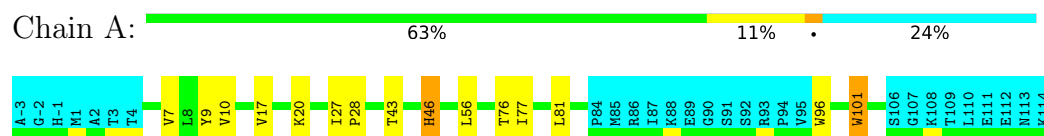


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.18 Score per residue for model 18

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

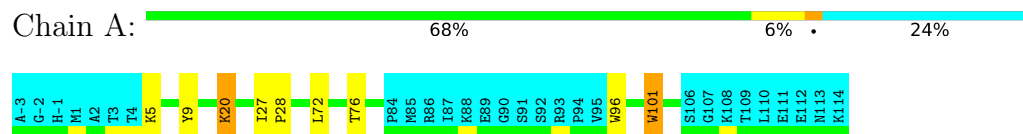


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.19 Score per residue for model 19

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E

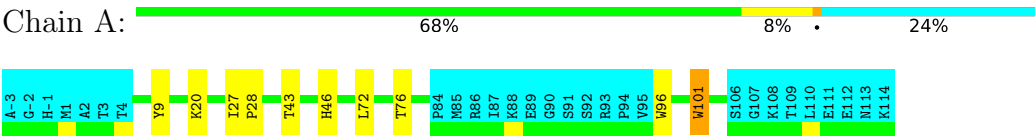


- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



4.2.20 Score per residue for model 20

- Molecule 1: Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: RNA (5'-R(*AP*AP*UP*AP*AP*A)-3')



5 Refinement protocol and experimental data overview

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

Of the 250 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
Amber	refinement	10
CYANA	structure calculation	3.96

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	1398
Number of shifts mapped to atoms	1323
Number of unparsed shifts	0
Number of shifts with mapping errors	75
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	78%

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	0.51±0.00	0±0/737 (0.0± 0.0%)	1.02±0.02	1±1/999 (0.1± 0.1%)
2	B	0.75±0.01	0±0/143 (0.0± 0.0%)	1.04±0.06	0±0/221 (0.1± 0.2%)
All	All	0.56	0/17600 (0.0%)	1.02	20/24400 (0.1%)

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
2	B	0.0±0.0	0.1±0.3
All	All	0	2

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	3	U	C2'-C3'-O3'	-6.31	104.23	113.70	16	4
1	A	46	HIS	CB-CG-CD2	-5.70	123.80	131.20	14	13
1	A	83	LYS	CA-C-N	5.21	124.93	119.82	2	1
1	A	83	LYS	C-N-CA	5.21	124.93	119.82	2	1
2	B	5	A	O4'-C1'-N9	5.02	115.73	108.20	12	1

There are no chirality outliers.

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
2	B	3	U	Sidechain	2

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	720	692	692	4±1
2	B	127	65	67	0±1
All	All	16940	15140	15180	82

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 3.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:101:TRP:CD1	1:A:101:TRP:C	0.54	2.85	5	20
1:A:20:LYS:H	1:A:20:LYS:CD	0.51	2.19	19	6
1:A:27:ILE:N	1:A:28:PRO:CD	0.49	2.76	18	20
2:B:3:U:C5'	2:B:3:U:C6	0.49	2.95	19	1
1:A:9:TYR:CD1	1:A:9:TYR:C	0.47	2.92	16	16
1:A:81:LEU:HB2	1:A:96:TRP:CZ2	0.47	2.44	6	3
2:B:3:U:C6	2:B:3:U:H5''	0.46	2.45	19	1
1:A:20:LYS:HD3	1:A:20:LYS:H	0.46	1.70	1	4
1:A:103:LYS:HA	1:A:103:LYS:HE3	0.45	1.89	6	1
1:A:81:LEU:HB2	1:A:96:TRP:CH2	0.43	2.47	13	1
1:A:20:LYS:H	1:A:20:LYS:HD3	0.42	1.74	19	1
2:B:3:U:C6	2:B:3:U:O5'	0.42	2.72	12	5
1:A:20:LYS:HD2	1:A:20:LYS:H	0.42	1.73	11	1
1:A:98:ASP:HA	1:A:101:TRP:CE3	0.41	2.50	5	1
2:B:3:U:O5'	2:B:3:U:C6	0.40	2.74	8	1

6.3 Torsion angles

6.3.1 Protein backbone

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	89/117 (76%)	84±2 (94±2%)	5±1 (5±2%)	0±0 (0±1%)	31	76
All	All	1780/2340 (76%)	1681 (94%)	91 (5%)	8 (0%)	31	76

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

Mol	Chain	Res	Type	Models (Total)
1	A	7	VAL	7
1	A	5	LYS	1

6.3.2 Protein sidechains ⓘ

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	74/97 (76%)	67±1 (91±2%)	7±1 (9±2%)	11	56
All	All	1480/1940 (76%)	1343 (91%)	137 (9%)	11	56

All 16 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	20	LYS	20
1	A	76	THR	20
1	A	101	TRP	20
1	A	43	THR	19
1	A	72	LEU	11
1	A	17	VAL	10
1	A	45	LYS	10
1	A	56	LEU	9
1	A	10	VAL	6
1	A	47	ARG	3
1	A	46	HIS	2
1	A	103	LYS	2
1	A	104	LYS	2
1	A	81	LEU	1
1	A	80	ASN	1
1	A	77	ILE	1

6.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers	Suiteness
2	B	5/6 (83%)	4±0 (80±0%)	1±1 (29±18%)	0.00±0.00
All	All	102/120 (85%)	80 (78%)	29 (28%)	0.00

The overall RNA backbone suiteness is 0.00.

All unique RNA backbone outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
2	B	2	A	20
2	B	3	U	20
2	B	4	A	20
2	B	5	A	20

All unique RNA pucker outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
2	B	3	U	14
2	B	5	A	13
2	B	1	A	2

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 oligosaccharides 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

The completeness of assignment taking into account all chemical shift lists is 78% for the well-defined parts and 78% for the entire structure.

7.1 Chemical shift list 1

File name: `working_cs.cif`

Chemical shift list name: *starch_output*

7.1.1 Bookkeeping

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

Total number of shifts	1398
Number of shifts mapped to atoms	1323
Number of unparsed shifts	0
Number of shifts with mapping errors	75
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. All 75 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	B	1	ADE	H1'	5.949	0.000	.
1	B	1	ADE	H2	8.062	0.000	1
1	B	1	ADE	H2'	4.8	0.000	.
1	B	1	ADE	H3'	4.744	0.000	.
1	B	1	ADE	H4'	4.417	0.000	.
1	B	1	ADE	H5"	3.855	0.000	.
1	B	1	ADE	H5'	3.873	0.000	.
1	B	1	ADE	H8	8.252	0.000	1
1	B	1	ADE	C1'	91.677	0.000	.
1	B	1	ADE	C2	154.993	0.000	1
1	B	1	ADE	C2'	75.701	0.000	.
1	B	1	ADE	C3'	76.878	0.000	.
1	B	1	ADE	C4'	86.95	0.000	.
1	B	1	ADE	C5'	63.834	0.000	.

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	B	1	ADE	C8	141.841	0.000	1
1	B	2	ADE	H1'	5.971	0.006	.
1	B	2	ADE	H2	8.084	0.000	1
1	B	2	ADE	H2'	4.783	0.000	.
1	B	2	ADE	H3'	4.849	0.000	.
1	B	2	ADE	H4'	4.567	0.000	.
1	B	2	ADE	H5"	4.252	0.000	.
1	B	2	ADE	H5'	4.362	0.000	.
1	B	2	ADE	H8	8.288	0.000	1
1	B	2	ADE	C1'	89.977	0.004	.
1	B	2	ADE	C2	155.374	0.000	1
1	B	2	ADE	C2'	76.195	0.000	.
1	B	2	ADE	C3'	76.668	0.000	.
1	B	2	ADE	C4'	85.231	0.000	.
1	B	2	ADE	C5'	67.441	0.013	.
1	B	2	ADE	C8	141.716	0.000	1
1	B	4	ADE	H1'	5.87	0.000	.
1	B	4	ADE	H2	8.03	0.000	1
1	B	4	ADE	H2'	4.746	0.000	.
1	B	4	ADE	H3'	4.758	0.000	.
1	B	4	ADE	H4'	4.526	0.000	.
1	B	4	ADE	H5"	4.193	0.000	.
1	B	4	ADE	H5'	4.284	0.000	.
1	B	4	ADE	H8	8.246	0.000	1
1	B	4	ADE	C1'	90.941	0.000	.
1	B	4	ADE	C2	155.178	0.000	1
1	B	4	ADE	C2'	76.099	0.000	.
1	B	4	ADE	C3'	76.386	0.000	.
1	B	4	ADE	C4'	84.718	0.000	.
1	B	4	ADE	C5'	67.65	0.012	.
1	B	4	ADE	C8	143.153	0.000	1
1	B	5	ADE	H1'	5.896	0.000	.
1	B	5	ADE	H2	7.909	0.000	1
1	B	5	ADE	H2'	4.641	0.000	.
1	B	5	ADE	H3'	4.813	0.000	.
1	B	5	ADE	H4'	4.527	0.000	.
1	B	5	ADE	H5"	4.233	0.000	.
1	B	5	ADE	H5'	4.383	0.000	.
1	B	5	ADE	H8	8.194	0.000	1
1	B	5	ADE	C1'	90.852	0.000	.
1	B	5	ADE	C2	154.878	0.000	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	B	5	ADE	C2'	76.402	0.000	.
1	B	5	ADE	C3'	75.483	0.000	.
1	B	5	ADE	C4'	84.398	0.000	.
1	B	5	ADE	C5'	66.979	0.000	.
1	B	5	ADE	C8	141.324	0.000	1
1	B	6	ADE	H1'	6.004	0.000	.
1	B	6	ADE	H2	8.131	0.000	1
1	B	6	ADE	H2'	4.535	0.000	.
1	B	6	ADE	H3'	4.502	0.000	.
1	B	6	ADE	H4'	4.392	0.000	.
1	B	6	ADE	H5"	4.217	0.000	.
1	B	6	ADE	H5'	4.381	0.000	.
1	B	6	ADE	H8	8.26	0.000	1
1	B	6	ADE	C1'	90.448	0.000	.
1	B	6	ADE	C2	155.413	0.000	1
1	B	6	ADE	C2'	77.399	0.000	.
1	B	6	ADE	C3'	72.254	0.000	.
1	B	6	ADE	C4'	85.762	0.000	.
1	B	6	ADE	C5'	67.278	0.029	.
1	B	6	ADE	C8	141.726	0.000	1

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	114	-0.17 ± 0.13	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	105	0.04 ± 0.09	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	106	0.43 ± 0.28	None needed (< 0.5 ppm)

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 78%, i.e. 1043 atoms were assigned a chemical shift out of a possible 1341. 0 out of 14 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	351/446 (79%)	178/181 (98%)	88/178 (49%)	85/87 (98%)

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	Total	¹H	¹³C	¹⁵N
Sidechain	583/657 (89%)	399/425 (94%)	180/210 (86%)	4/22 (18%)
Aromatic	96/126 (76%)	57/63 (90%)	37/59 (63%)	2/4 (50%)
Sugar	11/66 (17%)	6/36 (17%)	5/30 (17%)	0/0 (—%)
Base	2/46 (4%)	1/28 (4%)	1/12 (8%)	0/6 (0%)
Overall	1043/1341 (78%)	641/733 (87%)	311/489 (64%)	91/119 (76%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 78%, i.e. 1321 atoms were assigned a chemical shift out of a possible 1702. 0 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	448/585 (77%)	228/238 (96%)	114/234 (49%)	106/113 (94%)
Sidechain	760/872 (87%)	520/564 (92%)	235/276 (85%)	5/32 (16%)
Aromatic	100/133 (75%)	59/67 (88%)	39/61 (64%)	2/5 (40%)
Sugar	11/66 (17%)	6/36 (17%)	5/30 (17%)	0/0 (—%)
Base	2/46 (4%)	1/28 (4%)	1/12 (8%)	0/6 (0%)
Overall	1321/1702 (78%)	814/933 (87%)	394/613 (64%)	113/156 (72%)

7.1.4 Statistically unusual chemical shifts ⓘ

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots ⓘ

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:

