



Full wwPDB NMR Structure Validation Report i

Apr 20, 2024 – 03:40 PM EDT

PDB ID : 2N6T
BMRB ID : 25781
Title : NMR Assignment and NMR Structure of CssA3 (top stem) of CssA thermometer
Authors : Barnwal, R.; Godin, K.; Varani, G.; Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2015-08-28

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 types of validation reports are described at
<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

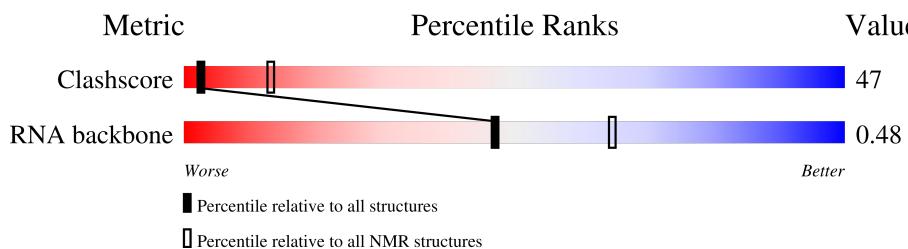
MolProbitiy : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

1 Overall quality at a glance

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

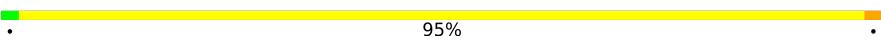
The overall completeness of chemical shifts assignment is 37%.

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
RNA backbone	4643	676

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

Mol	Chain	Length	Quality of chain	
1	A	42	 .	95%

2 Ensemble composition and analysis

This entry contains 10 models. This entry does not contain polypeptide chains, therefore identification of well-defined residues and clustering analysis are not possible. All residues are included in the validation scores.

3 Entry composition [\(i\)](#)

There is only 1 type of molecule in this entry. The entry contains 1329 atoms, of which 448 are hydrogens and 0 are deuteriums.

- Molecule 1 is a RNA chain called RNA (42-MER).

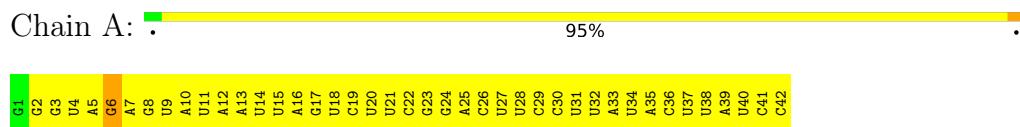
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
1	A	42	1329	396	448	146	298	41	0

4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

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

- Molecule 1: RNA (42-MER)

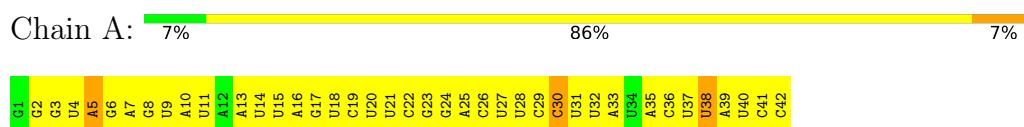


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

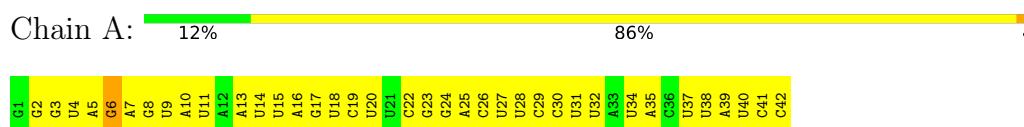
4.2.1 Score per residue for model 1

- Molecule 1: RNA (42-MER)



4.2.2 Score per residue for model 2

- Molecule 1: RNA (42-MER)



4.2.3 Score per residue for model 3

- Molecule 1: RNA (42-MER)



G1	G2	G3	G4	G5	G6	G7	G8	G9	G10	G11	G12	G13	G14	G15	G16	G17	G18	G19	G20	G21	G22	G23	G24	G25	G26	G27	G28	G29	G30	G31	G32	G33	G34	G35	G36	G37	G38	G39	G40	G41	G42
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4.2.4 Score per residue for model 4

- Molecule 1: RNA (42-MER)



G1	G2	G3	G4	G5	G6	G7	G8	G9	G10	G11	G12	G13	G14	G15	G16	G17	G18	G19	G20	G21	G22	G23	G24	G25	G26	G27	G28	G29	G30	G31	G32	G33	G34	G35	G36	G37	G38	G39
----	----	----	----	----	----	----	----	----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4.2.5 Score per residue for model 5

- Molecule 1: RNA (42-MER)



G1	G2	G3	G4	G5	G6	A7	A8	U9	A10	A11	A12	A13	A14	A15	A16	G17	U18	C19	U21	C22	G23	C24	A25	C26	U27	A33	U34	A35	C36	U37	U38	A39	A40	C41	C42
----	----	----	----	----	----	----	----	----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4.2.6 Score per residue for model 6

- Molecule 1: RNA (42-MER)



G1	G2	G3	G4	A5	G6	A7	G8	U9	A10	U11	A12	A13	U14	U15	A16	G17	U18	C19	U20	U21	C22	G23	C24	A25	C26	U27	U28	C29	C30	U31	U32	A33	A34	C35	C36	U37	U38	A39	A40	C41	C42
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4.2.7 Score per residue for model 7

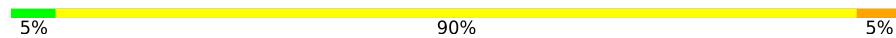
- Molecule 1: RNA (42-MER)

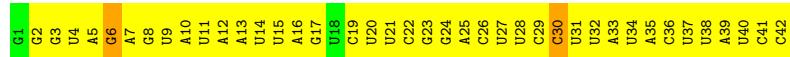


G1	G2	G3	U4	A5	G6	A7	U9	A10	U11	A12	A13	U14	U15	A16	G17	U18	C19	U20	U21	C22	G23	G24	A25	C26	U27	U28	C29	C30	U31	U32	A33	U34	A35	C36	U37	U38	A39	U40	C41	C42
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4.2.8 Score per residue for model 8

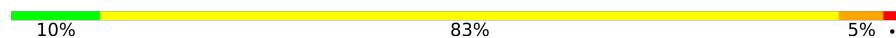
- Molecule 1: RNA (42-MER)

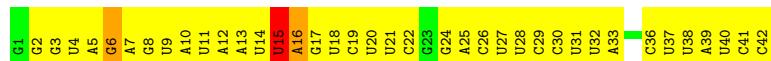
Chain A:  5% 5%



4.2.9 Score per residue for model 9

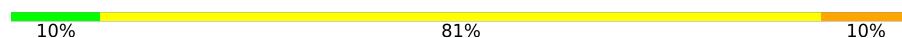
- Molecule 1: RNA (42-MER)

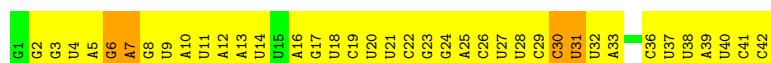
Chain A:  10% 83% 5%



4.2.10 Score per residue for model 10

- Molecule 1: RNA (42-MER)

Chain A:  10% 81% 10%



5 Refinement protocol and experimental data overview i

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

Of the 50 calculated structures, 10 were deposited, based on the following criterion: *structures with the least restraint violations*.

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

Software name	Classification	Version
X-PLOR NIH	structure solution	
X-PLOR NIH	refinement	

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	314
Number of shifts mapped to atoms	314
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	37%

6 Model quality [\(i\)](#)

6.1 Standard geometry [\(i\)](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	1.09±0.02	0±0/982 (0.0± 0.0%)	1.18±0.01	0±0/1526 (0.0± 0.0%)
All	All	1.09	0/9820 (0.0%)	1.18	1/15260 (0.0%)

There are no bond-length outliers.

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	15	U	C3'-C2'-C1'	5.34	105.78	101.50	9	1

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [\(i\)](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	881	448	448	63±8
All	All	8810	4480	4480	629

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 47.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:37:U:HO2'	1:A:38:U:H6	0.93	0.98	9	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:14:U:HO2'	1:A:15:U:H6	0.88	1.07	6	1
1:A:6:G:HO2'	1:A:7:A:H8	0.87	0.92	3	2
1:A:15:U:O2'	1:A:16:A:O4'	0.83	1.96	9	1
1:A:13:A:H2'	1:A:14:U:O4'	0.82	1.74	1	10
1:A:29:C:O2'	1:A:30:C:H5'	0.82	1.72	3	9
1:A:5:A:O2'	1:A:6:G:H5'	0.81	1.75	3	3
1:A:4:U:O2'	1:A:5:A:H5'	0.81	1.75	4	8
1:A:27:U:H2'	1:A:28:U:O4'	0.79	1.78	7	9
1:A:31:U:O2'	1:A:32:U:H5'	0.79	1.78	1	8
1:A:2:G:H2'	1:A:3:G:O4'	0.78	1.78	10	10
1:A:7:A:O2'	1:A:8:G:H5'	0.78	1.79	8	8
1:A:16:A:H2'	1:A:17:G:O4'	0.77	1.79	10	9
1:A:38:U:O2'	1:A:39:A:H5'	0.77	1.80	10	1
1:A:24:G:H2'	1:A:25:A:O4'	0.77	1.79	5	7
1:A:5:A:H2'	1:A:6:G:O4'	0.77	1.79	7	3
1:A:9:U:O2'	1:A:10:A:H5'	0.77	1.80	3	10
1:A:41:C:O2'	1:A:42:C:H5'	0.76	1.81	7	10
1:A:19:C:H2'	1:A:20:U:O4'	0.76	1.81	1	9
1:A:27:U:O2'	1:A:28:U:H5'	0.75	1.81	7	10
1:A:32:U:O2'	1:A:33:A:H5'	0.74	1.82	3	8
1:A:7:A:H2'	1:A:8:G:O4'	0.73	1.83	3	3
1:A:28:U:O2'	1:A:29:C:H5'	0.72	1.84	10	6
1:A:19:C:O2'	1:A:20:U:H5'	0.72	1.84	6	5
1:A:39:A:H2'	1:A:40:U:O4'	0.72	1.84	4	8
1:A:21:U:O2'	1:A:22:C:H5'	0.71	1.85	9	3
1:A:25:A:H2'	1:A:26:C:O4'	0.71	1.84	8	3
1:A:10:A:O2'	1:A:11:U:H5'	0.70	1.86	6	10
1:A:23:G:H2'	1:A:24:G:O4'	0.70	1.85	7	5
1:A:2:G:O2'	1:A:3:G:H5'	0.70	1.86	2	10
1:A:18:U:O2'	1:A:19:C:H5'	0.70	1.85	1	3
1:A:15:U:O2'	1:A:16:A:C8	0.68	2.42	9	1
1:A:16:A:O2'	1:A:17:G:H5'	0.68	1.89	5	8
1:A:39:A:O2'	1:A:40:U:H5'	0.67	1.90	1	10
1:A:23:G:O2'	1:A:24:G:H5'	0.67	1.90	8	3
1:A:6:G:O2'	1:A:7:A:H8	0.66	1.71	3	2
1:A:1:G:O2'	1:A:2:G:H5'	0.66	1.90	7	1
1:A:15:U:O2'	1:A:16:A:O5'	0.65	2.15	9	1
1:A:13:A:O2'	1:A:14:U:H5'	0.64	1.93	7	8
1:A:25:A:O2'	1:A:26:C:H5'	0.64	1.92	6	9
1:A:24:G:O2'	1:A:25:A:H5'	0.64	1.93	9	8
1:A:12:A:O2'	1:A:13:A:H5'	0.63	1.92	5	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:26:C:O2'	1:A:27:U:H5'	0.63	1.94	5	4
1:A:14:U:H1'	1:A:31:U:O4	0.63	1.94	9	2
1:A:37:U:O2'	1:A:38:U:H6	0.63	1.72	9	1
1:A:40:U:H6	1:A:40:U:O5'	0.62	1.77	4	8
1:A:38:U:O2'	1:A:39:A:O4'	0.62	2.17	1	1
1:A:14:U:O2	1:A:15:U:H5	0.61	1.78	8	1
1:A:37:U:H2'	1:A:38:U:O4'	0.61	1.95	2	2
1:A:28:U:H2'	1:A:29:C:O4'	0.60	1.95	10	7
1:A:25:A:O2'	1:A:26:C:O4'	0.60	2.18	5	6
1:A:5:A:O2'	1:A:6:G:C8	0.60	2.54	10	3
1:A:24:G:H8	1:A:24:G:O5'	0.59	1.79	6	10
1:A:32:U:H6	1:A:32:U:O5'	0.59	1.80	9	2
1:A:25:A:O2'	1:A:26:C:C5'	0.59	2.50	3	9
1:A:5:A:C6	1:A:39:A:C2	0.59	2.91	7	3
1:A:6:G:O2'	1:A:7:A:C8	0.58	2.56	5	4
1:A:35:A:O2'	1:A:36:C:H5'	0.58	1.98	1	3
1:A:5:A:C2	1:A:39:A:N1	0.58	2.71	9	4
1:A:5:A:N3	1:A:6:G:N7	0.57	2.52	1	1
1:A:5:A:C2	1:A:39:A:C2	0.56	2.93	10	2
1:A:19:C:C5	1:A:20:U:C4	0.56	2.93	9	4
1:A:7:A:N6	1:A:38:U:C6	0.56	2.74	9	1
1:A:5:A:N7	1:A:6:G:C2	0.55	2.74	5	1
1:A:3:G:O2'	1:A:4:U:H5'	0.55	2.01	10	1
1:A:21:U:O2'	1:A:23:G:C5	0.55	2.58	1	4
1:A:17:G:N2	1:A:27:U:C2	0.55	2.75	6	2
1:A:5:A:N6	1:A:38:U:C4	0.55	2.75	7	2
1:A:37:U:O2'	1:A:38:U:C6	0.54	2.61	10	1
1:A:5:A:N1	1:A:6:G:N2	0.54	2.55	10	1
1:A:4:U:O2'	1:A:5:A:C5'	0.54	2.56	1	7
1:A:22:C:H6	1:A:22:C:O5'	0.54	1.86	8	8
1:A:27:U:C4	1:A:28:U:C4	0.54	2.96	1	1
1:A:39:A:O2'	1:A:40:U:C5'	0.54	2.56	6	6
1:A:5:A:O2'	1:A:6:G:P	0.53	2.66	1	6
1:A:15:U:C4	1:A:16:A:C6	0.53	2.96	2	1
1:A:5:A:C6	1:A:6:G:N2	0.53	2.76	10	2
1:A:15:U:O2'	1:A:30:C:N4	0.53	2.41	1	1
1:A:13:A:C8	1:A:14:U:C6	0.53	2.97	8	2
1:A:28:U:C2	1:A:29:C:C2	0.53	2.97	5	3
1:A:27:U:C4	1:A:28:U:N3	0.53	2.77	1	1
1:A:22:C:O4'	1:A:23:G:C2	0.52	2.62	8	7
1:A:14:U:O2'	1:A:15:U:H6	0.52	1.87	8	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:5:A:N1	1:A:39:A:C2	0.52	2.76	3	2
1:A:38:U:O2'	1:A:39:A:C5'	0.52	2.56	10	2
1:A:41:C:O2'	1:A:42:C:C5'	0.52	2.58	3	5
1:A:8:G:O2'	1:A:9:U:H5'	0.52	2.05	4	6
1:A:38:U:O2'	1:A:39:A:O5'	0.52	2.28	1	1
1:A:27:U:O2'	1:A:28:U:C5'	0.52	2.56	5	8
1:A:7:A:N6	1:A:38:U:C5	0.52	2.78	6	1
1:A:25:A:C2	1:A:26:C:C2	0.52	2.98	7	3
1:A:28:U:HO2'	1:A:29:C:H6	0.51	1.48	1	2
1:A:5:A:N1	1:A:39:A:C6	0.51	2.79	9	2
1:A:34:U:O2'	1:A:35:A:H5'	0.51	2.05	8	6
1:A:27:U:C5	1:A:28:U:C4	0.51	2.99	1	1
1:A:37:U:C2	1:A:38:U:C4	0.51	2.99	1	1
1:A:6:G:O2'	1:A:7:A:H5'	0.51	2.06	9	1
1:A:3:G:O2'	1:A:4:U:C5'	0.51	2.59	10	1
1:A:39:A:O5'	1:A:39:A:H8	0.50	1.89	1	1
1:A:12:A:C6	1:A:33:A:N1	0.50	2.79	8	2
1:A:35:A:H2'	1:A:36:C:O4'	0.50	2.06	1	1
1:A:30:C:O2	1:A:31:U:H5	0.50	1.89	2	1
1:A:28:U:O2'	1:A:29:C:C5'	0.50	2.58	10	3
1:A:37:U:H3'	1:A:37:U:O2	0.50	2.06	4	1
1:A:37:U:C2	1:A:38:U:C5	0.50	3.00	1	1
1:A:5:A:HO2'	1:A:6:G:H5'	0.49	1.67	3	1
1:A:20:U:C2	1:A:22:C:C2	0.49	3.01	7	5
1:A:5:A:O2'	1:A:6:G:O5'	0.49	2.30	4	3
1:A:37:U:O2'	1:A:38:U:H5'	0.49	2.06	4	1
1:A:29:C:O2'	1:A:30:C:C5'	0.49	2.60	4	6
1:A:6:G:H8	1:A:6:G:O5'	0.49	1.90	3	2
1:A:5:A:C2	1:A:38:U:O4	0.49	2.66	9	2
1:A:19:C:C4	1:A:20:U:C4	0.49	3.00	9	1
1:A:31:U:O2'	1:A:32:U:C5'	0.49	2.61	2	3
1:A:36:C:C5	1:A:37:U:C4	0.49	3.01	5	1
1:A:31:U:O2'	1:A:32:U:O4'	0.49	2.30	6	2
1:A:6:G:N2	1:A:36:C:N4	0.49	2.60	10	1
1:A:20:U:N3	1:A:22:C:O2	0.48	2.46	2	6
1:A:14:U:O2	1:A:15:U:C5	0.48	2.62	8	2
1:A:39:A:O5'	1:A:39:A:C8	0.48	2.65	10	1
1:A:5:A:N1	1:A:37:U:O4	0.48	2.46	10	1
1:A:35:A:O2'	1:A:36:C:C5'	0.48	2.61	6	3
1:A:6:G:O6	1:A:39:A:C4	0.48	2.66	1	1
1:A:36:C:N3	1:A:37:U:O2	0.48	2.46	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:37:U:O2	1:A:37:U:C3'	0.48	2.61	4	2
1:A:5:A:N1	1:A:39:A:N6	0.48	2.62	9	1
1:A:28:U:N3	1:A:29:C:C2	0.48	2.82	5	2
1:A:17:G:N2	1:A:18:U:C2	0.48	2.82	7	3
1:A:20:U:O2'	1:A:21:U:H2'	0.48	2.09	9	1
1:A:5:A:O2'	1:A:6:G:C5'	0.47	2.57	3	1
1:A:10:A:HO2'	1:A:11:U:H5'	0.47	1.69	8	1
1:A:16:A:H8	1:A:16:A:O5'	0.47	1.93	6	2
1:A:15:U:O2'	1:A:16:A:P	0.47	2.72	9	1
1:A:24:G:O2'	1:A:25:A:C5'	0.47	2.63	4	1
1:A:7:A:N6	1:A:38:U:O2	0.47	2.47	5	4
1:A:10:A:O2'	1:A:11:U:C5'	0.47	2.62	7	5
1:A:16:A:O2'	1:A:17:G:C5'	0.47	2.62	1	3
1:A:29:C:O2	1:A:30:C:C6	0.47	2.68	1	1
1:A:9:U:O2'	1:A:10:A:C5'	0.47	2.63	10	3
1:A:36:C:O2'	1:A:37:U:H5'	0.47	2.09	3	2
1:A:37:U:O2	1:A:38:U:C5	0.47	2.68	10	2
1:A:7:A:C6	1:A:38:U:C6	0.47	3.03	9	1
1:A:2:G:O2'	1:A:3:G:C5'	0.46	2.62	1	2
1:A:29:C:N4	1:A:30:C:N4	0.46	2.63	8	1
1:A:36:C:H2'	1:A:37:U:O4'	0.46	2.10	1	4
1:A:32:U:H2'	1:A:33:A:O4'	0.46	2.11	10	1
1:A:8:G:N3	1:A:38:U:O4	0.46	2.48	10	1
1:A:26:C:O2'	1:A:27:U:C5'	0.46	2.63	4	2
1:A:7:A:N1	1:A:38:U:O2	0.46	2.49	5	2
1:A:39:A:C2'	1:A:40:U:O4'	0.46	2.64	3	4
1:A:7:A:N7	1:A:38:U:N3	0.46	2.64	9	1
1:A:28:U:O2'	1:A:29:C:H6	0.45	1.93	1	2
1:A:5:A:N6	1:A:38:U:O4	0.45	2.49	3	1
1:A:19:C:C4	1:A:20:U:N3	0.45	2.83	10	1
1:A:6:G:O2'	1:A:7:A:N7	0.45	2.47	4	2
1:A:13:A:C5	1:A:14:U:C2	0.45	3.04	3	3
1:A:7:A:N7	1:A:38:U:C2	0.45	2.84	9	1
1:A:17:G:O2'	1:A:18:U:O4'	0.45	2.34	1	2
1:A:10:A:H2'	1:A:11:U:O4'	0.45	2.11	6	2
1:A:7:A:C6	1:A:38:U:O2	0.45	2.69	8	1
1:A:5:A:N6	1:A:37:U:O4	0.45	2.49	10	1
1:A:28:U:O2'	1:A:29:C:C6	0.45	2.69	2	1
1:A:8:G:O2'	1:A:9:U:O4'	0.45	2.35	7	2
1:A:4:U:O2'	1:A:5:A:O4'	0.44	2.36	3	1
1:A:6:G:O5'	1:A:6:G:C8	0.44	2.71	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:11:U:O2'	1:A:12:A:O4'	0.44	2.34	3	2
1:A:12:A:N6	1:A:33:A:C6	0.44	2.85	8	2
1:A:16:A:C2'	1:A:17:G:O4'	0.44	2.66	2	2
1:A:30:C:O2'	1:A:31:U:OP2	0.44	2.36	3	2
1:A:12:A:O2'	1:A:13:A:C5'	0.44	2.64	7	1
1:A:18:U:O2'	1:A:19:C:C5'	0.44	2.61	1	1
1:A:20:U:O2	1:A:22:C:C2	0.44	2.71	5	2
1:A:33:A:O2'	1:A:34:U:O4'	0.44	2.32	8	2
1:A:5:A:N1	1:A:38:U:O4	0.44	2.51	6	1
1:A:32:U:O2'	1:A:33:A:C5'	0.43	2.63	1	1
1:A:29:C:H2'	1:A:30:C:O4'	0.43	2.13	5	2
1:A:29:C:HO2'	1:A:30:C:H5'	0.43	1.71	3	1
1:A:37:U:O2	1:A:37:U:H3'	0.43	2.13	6	1
1:A:5:A:C2	1:A:39:A:C6	0.43	3.06	9	1
1:A:8:G:C2	1:A:37:U:O2	0.43	2.71	2	1
1:A:13:A:C8	1:A:14:U:C5	0.43	3.06	8	2
1:A:4:U:C4	1:A:5:A:C6	0.43	3.06	4	1
1:A:7:A:O2'	1:A:8:G:C5'	0.43	2.60	1	1
1:A:30:C:HO2'	1:A:31:U:P	0.43	2.36	5	1
1:A:7:A:HO2'	1:A:8:G:H5'	0.43	1.73	5	1
1:A:25:A:HO2'	1:A:26:C:C4'	0.43	2.24	7	2
1:A:17:G:O2'	1:A:18:U:H5'	0.43	2.13	5	1
1:A:8:G:N2	1:A:37:U:O2	0.43	2.51	10	1
1:A:3:G:O2'	1:A:4:U:O4'	0.42	2.35	10	1
1:A:36:C:O2'	1:A:37:U:O4'	0.42	2.38	1	1
1:A:20:U:C2	1:A:22:C:O2	0.42	2.73	3	1
1:A:6:G:O6	1:A:39:A:N3	0.42	2.52	1	1
1:A:25:A:C2'	1:A:26:C:O4'	0.42	2.68	1	1
1:A:7:A:C6	1:A:38:U:C2	0.42	3.07	7	1
1:A:17:G:N2	1:A:18:U:O2	0.42	2.52	7	1
1:A:21:U:O2'	1:A:23:G:N7	0.42	2.50	7	1
1:A:4:U:O4	1:A:5:A:N6	0.42	2.53	4	1
1:A:36:C:C4	1:A:37:U:N3	0.42	2.87	5	1
1:A:14:U:O2'	1:A:15:U:C6	0.42	2.70	8	1
1:A:26:C:O2	1:A:26:C:C2'	0.42	2.68	8	1
1:A:29:C:O2	1:A:30:C:C5	0.42	2.73	1	1
1:A:17:G:C2	1:A:18:U:C2	0.42	3.08	7	1
1:A:7:A:C6	1:A:8:G:C4	0.41	3.08	3	1
1:A:29:C:C4	1:A:30:C:N3	0.41	2.88	2	1
1:A:35:A:O2'	1:A:36:C:O4'	0.41	2.37	7	2
1:A:28:U:C2'	1:A:29:C:O4'	0.41	2.67	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:37:U:C2'	1:A:38:U:O4'	0.41	2.67	2	2
1:A:5:A:H62	1:A:6:G:N2	0.41	2.12	5	1
1:A:37:U:O2	1:A:38:U:H5	0.41	1.98	10	1
1:A:24:G:O5'	1:A:24:G:C8	0.41	2.69	6	1
1:A:7:A:O5'	1:A:7:A:H8	0.41	1.99	1	1
1:A:28:U:HO2'	1:A:29:C:C5'	0.41	2.29	1	1
1:A:36:C:C2	1:A:37:U:C2	0.41	3.08	1	1
1:A:15:U:H2'	1:A:16:A:C8	0.41	2.51	1	1
1:A:7:A:N6	1:A:38:U:N3	0.41	2.69	7	1
1:A:12:A:C6	1:A:33:A:C6	0.41	3.09	8	1
1:A:7:A:C8	1:A:38:U:C2	0.41	3.09	9	1
1:A:5:A:N6	1:A:6:G:N2	0.40	2.69	5	1
1:A:28:U:N3	1:A:29:C:O2	0.40	2.54	5	1
1:A:25:A:C2	1:A:26:C:O2	0.40	2.74	7	1
1:A:36:C:C2'	1:A:37:U:O4'	0.40	2.69	1	1
1:A:28:U:O2'	1:A:29:C:O4'	0.40	2.40	7	1

6.3 Torsion angles [\(i\)](#)

6.3.1 Protein backbone [\(i\)](#)

There are no protein molecules in this entry.

6.3.2 Protein sidechains [\(i\)](#)

There are no protein molecules in this entry.

6.3.3 RNA [\(i\)](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers	Suiteness
1	A	41/42 (98%)	2±1 (4±3%)	1±1 (2±2%)	0.48±0.03
All	All	410/420 (98%)	18 (4%)	7 (2%)	0.48

The overall RNA backbone suiteness is 0.48.

All unique RNA backbone outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
1	A	6	G	6

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Mol	Chain	Res	Type	Models (Total)
1	A	16	A	4
1	A	30	C	2
1	A	7	A	2
1	A	31	U	2
1	A	38	U	1
1	A	15	U	1

All unique RNA pucker outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
1	A	30	C	3
1	A	6	G	3
1	A	5	A	1

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

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

6.5 Carbohydrates [\(i\)](#)

There are no monosaccharides in this entry.

6.6 Ligand geometry [\(i\)](#)

There are no ligands in this entry.

6.7 Other polymers [\(i\)](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

7 Chemical shift validation (i)

The completeness of assignment taking into account all chemical shift lists is 37% for the well-defined parts and 37% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping (i)

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

Total number of shifts	314
Number of shifts mapped to atoms	314
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	2

7.1.2 Chemical shift referencing (i)

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

7.1.3 Completeness of resonance assignments (i)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 37%, i.e. 286 atoms were assigned a chemical shift out of a possible 774. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Sugar	128/462 (28%)	109/252 (43%)	19/210 (9%)	0/0 (—%)
Base	158/312 (51%)	105/186 (56%)	39/76 (51%)	14/50 (28%)
Overall	286/774 (37%)	214/438 (49%)	58/286 (20%)	14/50 (28%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 37%, i.e. 286 atoms were assigned a chemical shift out of a possible 774. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Sugar	128/462 (28%)	109/252 (43%)	19/210 (9%)	0/0 (—%)
Base	158/312 (51%)	105/186 (56%)	39/76 (51%)	14/50 (28%)
Overall	286/774 (37%)	214/438 (49%)	58/286 (20%)	14/50 (28%)

7.1.4 Statistically unusual chemical shifts [\(i\)](#)

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

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	27	U	C6	0.00	131.00 – 152.02	-67.3
1	A	7	A	C1'	0.00	82.09 – 100.58	-49.4

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

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