



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 04:43 PM BST

PDB ID : 1QFQ
Title : Bacteriophage Lambda N-protein-NutboxB-RNA Complex
Authors : Schaerpf, M.; Sticht, H.; Roesch, P.
Deposited on : 1999-04-12

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<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

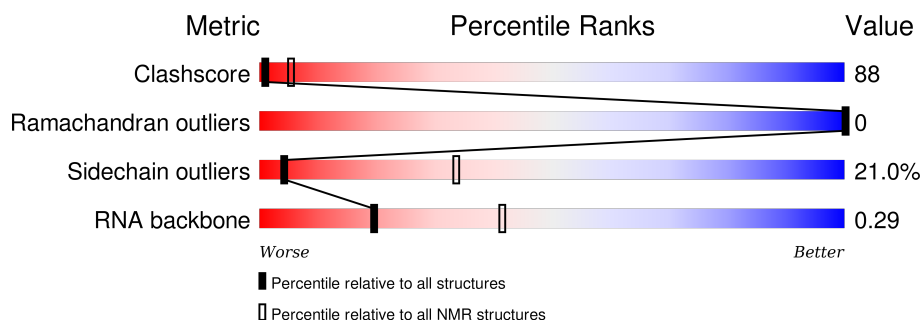
1 Overall quality at a glance

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
RNA backbone	3027	600

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	15	
2	B	35	

2 Ensemble composition and analysis

This entry contains 29 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative.

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	B:7-B:23 (17)	0.06	7

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 4 clusters and 2 single-model clusters were found.

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

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 1067 atoms, of which 460 are hydrogens and 0 are deuteriums.

- Molecule 1 is a RNA chain called 15-mer nutRboxB RNA hairpin.

Mol	Chain	Residues	Atoms						Trace
1	A	15	Total	C	H	N	O	P	0
			489	145	164	64	101	15	

- Molecule 2 is a protein called 36-mer N-terminal peptide of the N protein.

Mol	Chain	Residues	Atoms					Trace
2	B	35	Total	C	H	N	O	0
			578	171	296	62	49	

4 Residue-property plots

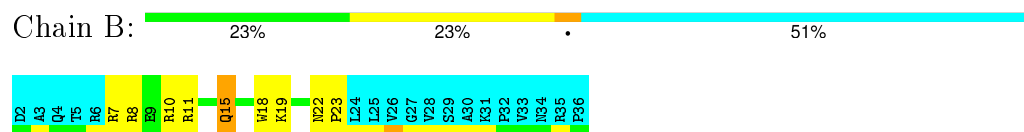
4.1 Average score per residue in the NMR ensemble

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: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein



4.2 Scores per residue for each member of the ensemble

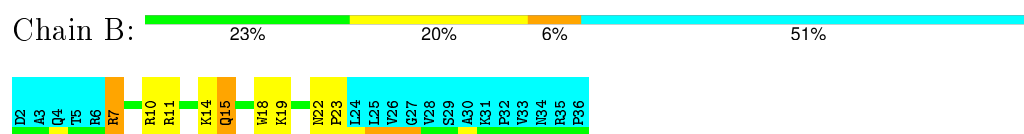
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

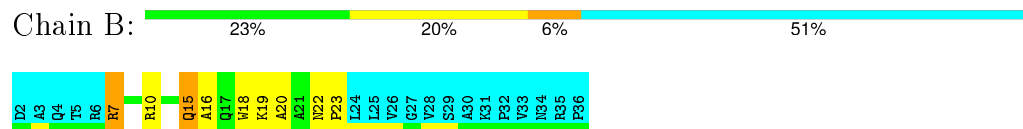


4.2.2 Score per residue for model 2

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

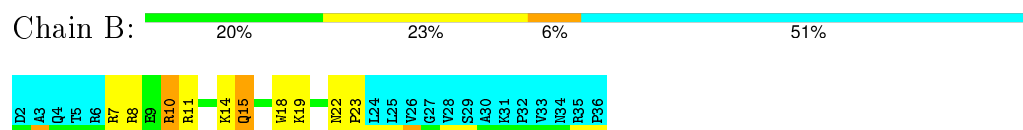


4.2.3 Score per residue for model 3

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

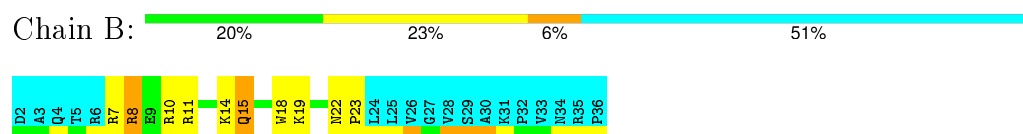


4.2.4 Score per residue for model 4

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

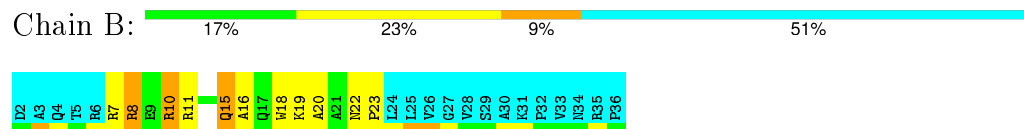


4.2.5 Score per residue for model 5

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

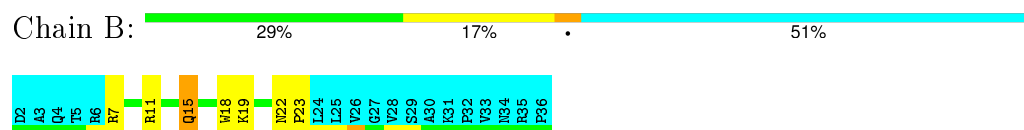


4.2.6 Score per residue for model 6

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

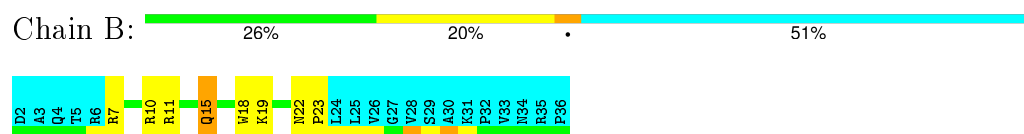


4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

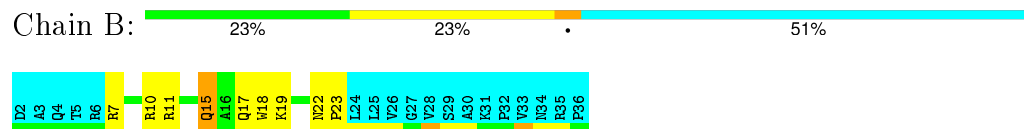


4.2.8 Score per residue for model 8

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

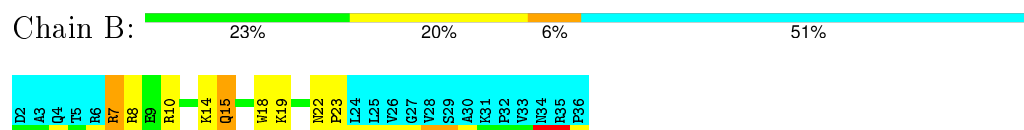


4.2.9 Score per residue for model 9

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

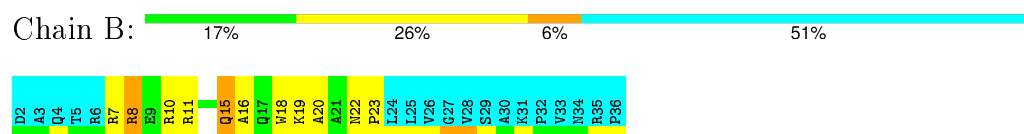


4.2.10 Score per residue for model 10

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

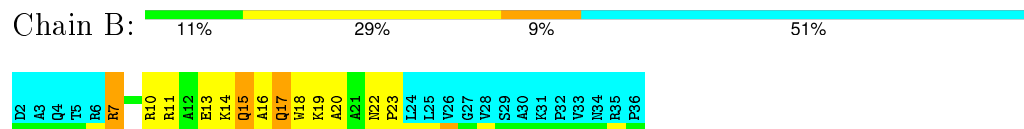


4.2.11 Score per residue for model 11

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

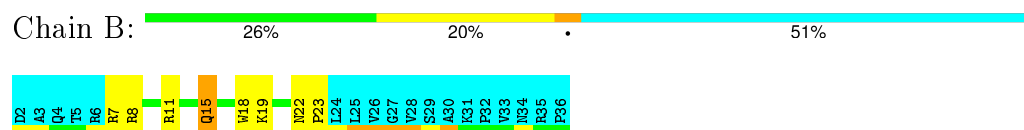


4.2.12 Score per residue for model 12

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

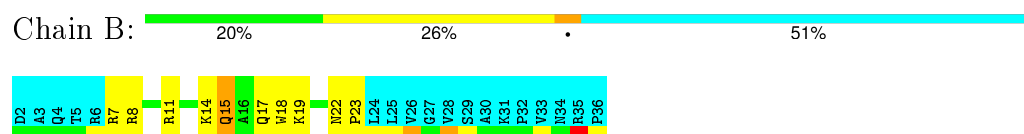


4.2.13 Score per residue for model 13

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

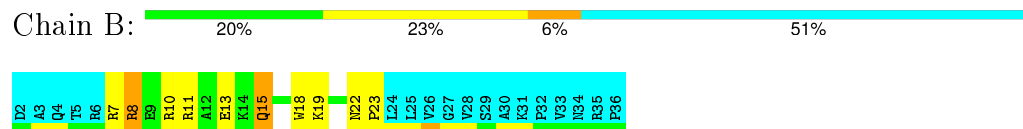


4.2.14 Score per residue for model 14

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

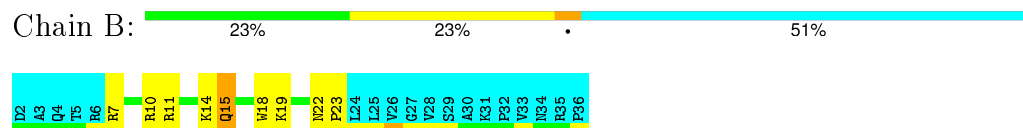


4.2.15 Score per residue for model 15

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

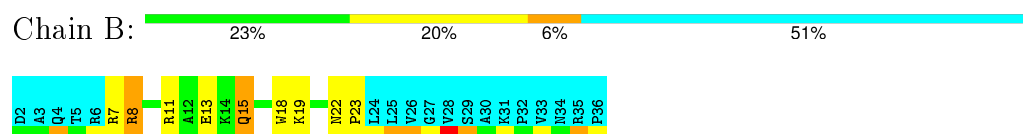


4.2.16 Score per residue for model 16

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

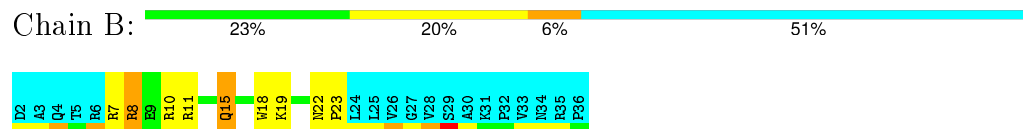


4.2.17 Score per residue for model 17

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

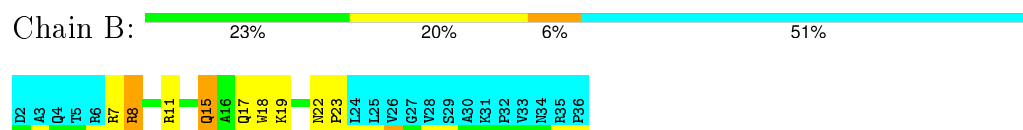


4.2.18 Score per residue for model 18

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

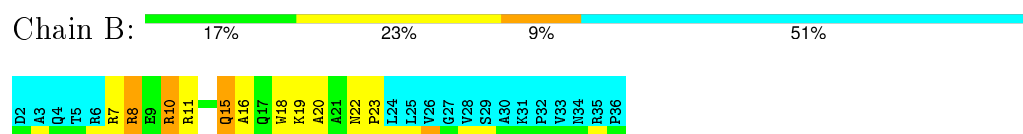


4.2.19 Score per residue for model 19

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

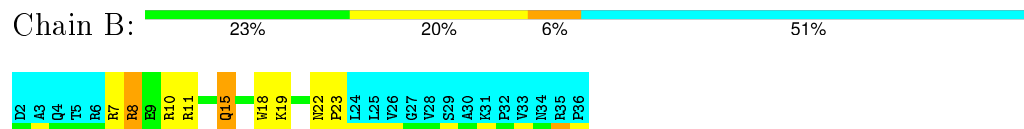


4.2.20 Score per residue for model 20

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

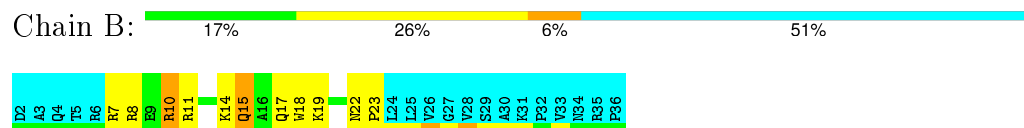


4.2.21 Score per residue for model 21

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

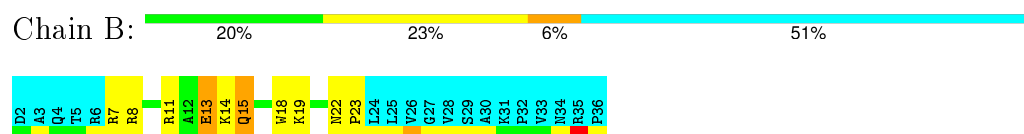


4.2.22 Score per residue for model 22

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

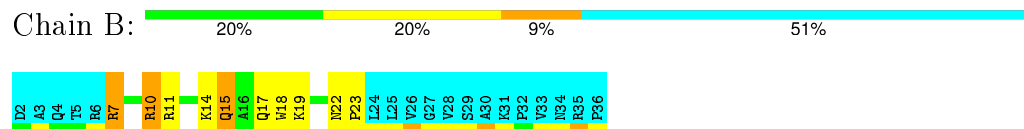


4.2.23 Score per residue for model 23

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

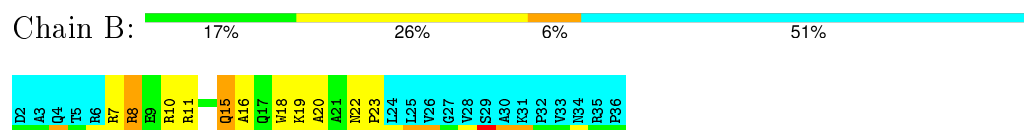


4.2.24 Score per residue for model 24

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

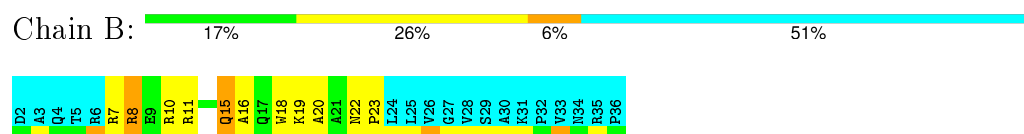


4.2.25 Score per residue for model 25

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

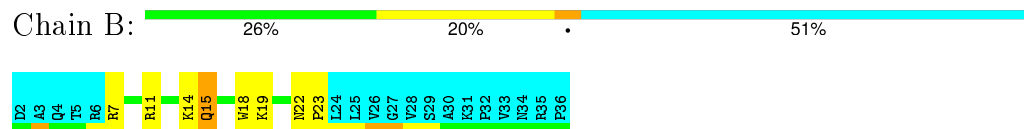


4.2.26 Score per residue for model 26

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

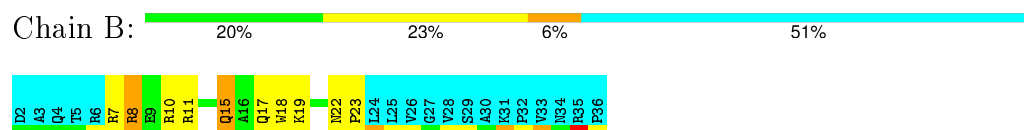


4.2.27 Score per residue for model 27

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein

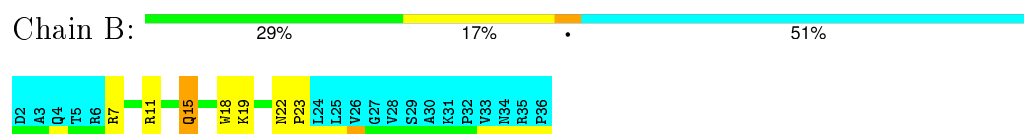


4.2.28 Score per residue for model 28

- Molecule 1: 15-mer nutRboxB RNA hairpin



- Molecule 2: 36-mer N-terminal peptide of the N protein



4.2.29 Score per residue for model 29

- Molecule 1: 15-mer nutRboxB RNA hairpin

Chain A:  13% 87%

G1	C2	C3	C4	U5	G6	A7	A8	A9	A10	A11	G12	G13	G14	C15
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- Molecule 2: 36-mer N-terminal peptide of the N protein

Chain B:  14% 26% 9% 51%

D2	A3	Q4	T5	R6	R7	R8	R9	R10	R11	K14	Q15	A16	Q17	M18	K19	A20	A21	N22	P23	L24	L25	V26	G27	V28	S29	A30	K31	P32	V33	N34	R35	P36
----	----	----	----	----	----	----	----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

5 Refinement protocol and experimental data overview

The models were refined using the following method: *MOLECULAR DYNAMICS*.

Of the 100 calculated structures, 29 were deposited, based on the following criterion: *LOWEST ENERGY, AGREEMENT WITH EXPERIMENTAL DATA*.

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

Software name	Classification	Version
X-PLOR	refinement	3.851
NDEE	structure solution	
XNDEE	structure solution	
X-PLOR	structure solution	

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

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	325	164	165	60±5
2	B	147	149	149	24±3
All	All	13688	9077	9106	2006

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 88.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:7:A:O2'	1:A:8:A:O5'	0.92	1.85	28	1
1:A:7:A:O4'	2:B:18:TRP:CG	0.87	2.27	21	29
1:A:11:A:H2'	1:A:12:G:O4'	0.85	1.72	5	29
1:A:5:U:H2'	1:A:6:G:H5'	0.81	1.52	26	20
1:A:11:A:O2'	1:A:12:G:OP1	0.81	1.99	6	22
1:A:12:G:OP1	1:A:12:G:H4'	0.80	1.75	25	10
1:A:5:U:C2'	1:A:6:G:H5'	0.80	2.07	26	16
1:A:12:G:H4'	1:A:12:G:OP1	0.78	1.78	1	12
1:A:3:C:O2'	1:A:4:C:OP1	0.78	2.00	2	22
1:A:3:C:O2'	1:A:4:C:P	0.77	2.43	27	22
1:A:12:G:O2'	1:A:13:G:OP1	0.76	2.02	7	11
1:A:1:G:O2'	1:A:2:C:P	0.74	2.46	26	17
1:A:13:G:O2'	1:A:14:G:P	0.73	2.45	24	16

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:12:G:O2'	1:A:13:G:P	0.73	2.47	7	7
1:A:14:G:O2'	1:A:15:C:P	0.72	2.47	11	2
1:A:12:G:H2'	1:A:13:G:C8	0.72	2.19	24	1
1:A:4:C:O2'	1:A:5:U:P	0.72	2.48	4	12
1:A:11:A:C6	1:A:12:G:C5	0.72	2.78	16	29
1:A:7:A:C8	2:B:18:TRP:NE1	0.70	2.59	28	29
1:A:1:G:O2'	1:A:2:C:OP1	0.69	2.09	12	17
1:A:4:C:N3	1:A:5:U:C5	0.69	2.61	18	29
1:A:4:C:C2	1:A:5:U:C6	0.68	2.82	2	24
1:A:12:G:C4'	1:A:12:G:OP1	0.68	2.42	9	12
1:A:4:C:O2'	1:A:5:U:OP1	0.68	2.12	2	12
1:A:12:G:OP1	1:A:12:G:C4'	0.68	2.42	19	10
1:A:12:G:C2'	1:A:13:G:OP1	0.67	2.42	15	7
1:A:7:A:H2'	1:A:8:A:C8	0.67	2.24	28	4
1:A:11:A:C5	1:A:12:G:C8	0.66	2.84	20	29
2:B:7:ARG:NH2	2:B:11:ARG:HD2	0.65	2.06	26	15
1:A:7:A:C8	2:B:18:TRP:CD1	0.65	2.84	28	28
1:A:1:G:C2'	1:A:2:C:O5'	0.65	2.44	19	4
1:A:14:G:O2'	1:A:15:C:OP1	0.65	2.13	11	2
1:A:6:G:N3	1:A:8:A:C8	0.65	2.65	26	29
1:A:7:A:C4	2:B:18:TRP:CE2	0.64	2.85	28	28
1:A:7:A:H1'	2:B:18:TRP:CD2	0.64	2.28	29	29
1:A:7:A:H4'	1:A:8:A:OP1	0.64	1.92	28	1
1:A:11:A:C2'	1:A:12:G:OP1	0.64	2.45	6	22
1:A:9:A:H61	2:B:19:LYS:NZ	0.64	1.89	14	2
1:A:6:G:N3	1:A:8:A:N7	0.64	2.46	9	25
1:A:8:A:OP1	2:B:15:GLN:HG3	0.64	1.92	28	1
1:A:7:A:H2'	1:A:8:A:O4'	0.63	1.93	29	29
1:A:8:A:N6	1:A:10:A:C6	0.63	2.67	15	26
1:A:7:A:HO2'	1:A:8:A:C5'	0.62	2.05	28	1
1:A:13:G:O2'	1:A:14:G:OP1	0.62	2.16	24	16
1:A:9:A:H61	2:B:19:LYS:CE	0.62	2.07	28	6
2:B:18:TRP:CD1	2:B:22:ASN:ND2	0.61	2.69	21	29
1:A:11:A:C6	1:A:12:G:N7	0.60	2.69	5	26
1:A:5:U:C2'	1:A:6:G:O5'	0.60	2.49	6	13
1:A:5:U:C4	1:A:6:G:N7	0.60	2.68	21	13
1:A:4:C:C2	1:A:5:U:C5	0.60	2.90	26	27
1:A:8:A:O2'	1:A:9:A:H3'	0.60	1.97	20	24
1:A:14:G:C6	1:A:15:C:N4	0.60	2.69	3	3
2:B:18:TRP:CZ2	2:B:22:ASN:HB3	0.60	2.31	29	29
1:A:7:A:C4	2:B:18:TRP:CZ2	0.60	2.90	28	27

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:14:G:C5	1:A:15:C:C5	0.60	2.90	8	15
1:A:7:A:O2'	2:B:15:GLN:CG	0.60	2.50	21	28
1:A:6:G:H4'	1:A:7:A:OP1	0.59	1.96	21	1
1:A:3:C:N3	1:A:4:C:C5	0.59	2.70	15	13
1:A:14:G:O2'	1:A:15:C:H5'	0.59	1.98	21	11
1:A:13:G:OP1	1:A:13:G:O4'	0.59	2.20	5	2
1:A:3:C:H2'	1:A:4:C:C6	0.58	2.33	13	17
1:A:13:G:H4'	1:A:14:G:OP1	0.58	1.98	26	10
1:A:7:A:C1'	2:B:18:TRP:CD2	0.58	2.87	21	29
1:A:8:A:OP1	2:B:8:ARG:HD3	0.58	1.99	21	5
1:A:7:A:O4'	2:B:18:TRP:CD1	0.57	2.57	21	1
1:A:6:G:H2'	1:A:7:A:H3'	0.57	1.75	28	2
1:A:6:G:O2'	1:A:7:A:P	0.57	2.62	21	1
1:A:3:C:C2	1:A:4:C:C5	0.57	2.92	7	21
1:A:3:C:C2	1:A:4:C:C6	0.57	2.92	12	7
1:A:4:C:O2'	1:A:5:U:H5'	0.57	2.00	1	2
1:A:13:G:H2'	1:A:14:G:C8	0.57	2.35	18	23
1:A:13:G:C2	1:A:14:G:C5	0.57	2.93	13	1
1:A:13:G:O4'	1:A:13:G:OP1	0.56	2.23	18	2
1:A:3:C:O2'	1:A:4:C:O5'	0.56	2.23	26	22
1:A:9:A:H61	2:B:19:LYS:HE2	0.56	1.59	12	4
1:A:13:G:N3	1:A:14:G:C8	0.56	2.74	13	1
2:B:22:ASN:N	2:B:23:PRO:CD	0.56	2.69	29	29
1:A:8:A:N6	1:A:10:A:N6	0.56	2.53	8	11
1:A:10:A:C2'	1:A:11:A:O5'	0.56	2.53	15	1
1:A:7:A:O2'	2:B:15:GLN:HG3	0.56	2.01	21	25
1:A:14:G:C2	1:A:15:C:C5	0.55	2.95	19	1
1:A:1:G:O2'	1:A:2:C:O5'	0.55	2.25	20	16
1:A:13:G:C2	1:A:14:G:N7	0.55	2.74	13	1
1:A:6:G:O6	2:B:8:ARG:CZ	0.55	2.54	25	12
1:A:11:A:C4	1:A:12:G:C8	0.55	2.95	12	13
1:A:9:A:H61	2:B:19:LYS:HZ3	0.55	1.42	14	1
1:A:6:G:N2	1:A:10:A:C8	0.55	2.75	21	28
1:A:14:G:C4	1:A:15:C:C5	0.55	2.95	19	3
1:A:3:C:H2'	1:A:4:C:H6	0.55	1.61	13	20
1:A:13:G:O2'	1:A:14:G:O5'	0.55	2.25	18	15
1:A:9:A:OP1	1:A:9:A:C4	0.55	2.60	21	10
1:A:5:U:H2'	1:A:6:G:O5'	0.54	2.02	5	13
1:A:3:C:C4	1:A:4:C:C5	0.54	2.95	26	3
1:A:3:C:N4	1:A:4:C:N4	0.54	2.54	29	3
1:A:7:A:HO2'	1:A:8:A:P	0.54	2.24	28	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:8:A:C6	1:A:10:A:N7	0.54	2.75	4	11
1:A:8:A:P	2:B:15:GLN:NE2	0.54	2.81	28	1
1:A:4:C:C4	1:A:5:U:C5	0.54	2.96	2	12
1:A:3:C:HO2'	1:A:4:C:P	0.54	2.17	23	3
1:A:14:G:HO2'	1:A:15:C:P	0.54	2.24	11	2
1:A:6:G:H1'	1:A:8:A:N7	0.54	2.18	24	27
1:A:2:C:H2'	1:A:3:C:O4'	0.54	2.03	1	6
1:A:14:G:C6	1:A:15:C:C4	0.53	2.96	8	10
1:A:7:A:O2'	1:A:8:A:P	0.53	2.66	28	1
1:A:9:A:P	2:B:8:ARG:HG2	0.53	2.42	24	1
1:A:8:A:P	2:B:15:GLN:HG3	0.53	2.43	28	1
1:A:7:A:O4'	2:B:18:TRP:CB	0.53	2.57	21	2
1:A:13:G:H2'	1:A:14:G:H8	0.53	1.63	2	10
1:A:10:A:HO2'	1:A:11:A:P	0.53	2.26	15	1
1:A:1:G:H2'	1:A:2:C:O5'	0.53	2.04	13	2
1:A:12:G:H2'	1:A:13:G:OP1	0.52	2.04	15	4
1:A:14:G:C6	1:A:15:C:C5	0.52	2.97	19	3
1:A:9:A:C4	1:A:9:A:OP1	0.52	2.63	28	18
1:A:9:A:N6	2:B:19:LYS:CE	0.52	2.73	22	5
1:A:9:A:H61	2:B:19:LYS:HE3	0.52	1.64	28	3
1:A:7:A:C1'	2:B:18:TRP:CG	0.52	2.93	24	29
1:A:8:A:O2'	1:A:9:A:C8	0.52	2.63	7	24
1:A:9:A:N6	2:B:19:LYS:HE3	0.52	2.20	28	3
1:A:5:U:H2'	1:A:6:G:C5'	0.52	2.35	16	12
1:A:9:A:N9	1:A:9:A:OP1	0.51	2.43	25	9
1:A:5:U:OP2	2:B:11:ARG:NH1	0.51	2.43	1	1
1:A:4:C:C2'	1:A:5:U:H5''	0.51	2.36	29	12
2:B:13:GLU:HG3	2:B:14:LYS:N	0.51	2.20	22	1
1:A:6:G:H3'	1:A:6:G:OP2	0.51	2.06	18	8
1:A:9:A:OP1	1:A:9:A:C8	0.51	2.64	16	11
1:A:10:A:H2'	1:A:11:A:O5'	0.51	2.05	15	1
1:A:5:U:OP2	2:B:7:ARG:NH2	0.51	2.43	6	9
2:B:22:ASN:N	2:B:23:PRO:HD3	0.51	2.21	29	14
1:A:14:G:N1	1:A:15:C:C4	0.51	2.78	19	2
1:A:8:A:C6	1:A:10:A:C5	0.51	2.99	27	25
1:A:9:A:N6	2:B:19:LYS:NZ	0.51	2.58	21	1
1:A:14:G:O2'	1:A:15:C:O5'	0.51	2.28	11	2
1:A:3:C:C4	1:A:4:C:C4	0.51	2.99	26	1
1:A:9:A:OP1	1:A:9:A:N9	0.50	2.45	20	15
2:B:7:ARG:CZ	2:B:11:ARG:NE	0.50	2.75	10	20
1:A:11:A:C5	1:A:12:G:N7	0.50	2.80	5	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:6:G:OP2	1:A:6:G:H3'	0.50	2.07	25	3
1:A:7:A:N9	2:B:18:TRP:CE2	0.49	2.80	28	20
1:A:8:A:OP1	2:B:15:GLN:CG	0.49	2.60	28	1
1:A:4:C:O2'	1:A:5:U:O5'	0.49	2.30	2	11
2:B:18:TRP:CE2	2:B:22:ASN:CB	0.49	2.96	29	17
2:B:18:TRP:CE3	2:B:19:LYS:N	0.49	2.81	6	29
1:A:4:C:H3'	2:B:7:ARG:NH1	0.49	2.23	4	7
1:A:8:A:O5'	2:B:15:GLN:HG3	0.48	2.08	28	1
1:A:5:U:O2'	1:A:6:G:P	0.48	2.71	6	11
1:A:7:A:C2'	1:A:8:A:O4'	0.48	2.61	28	1
1:A:11:A:C2	1:A:12:G:C4	0.48	3.02	21	3
1:A:4:C:H2'	1:A:5:U:O4'	0.48	2.08	2	3
2:B:18:TRP:CZ2	2:B:22:ASN:CB	0.48	2.97	29	6
1:A:3:C:O2'	1:A:4:C:H5'	0.48	2.09	1	1
1:A:9:A:H61	2:B:19:LYS:HD3	0.47	1.69	26	1
1:A:5:U:HO2'	1:A:6:G:P	0.47	2.31	16	10
1:A:10:A:O2'	1:A:11:A:P	0.47	2.72	15	1
2:B:16:ALA:O	2:B:20:ALA:HB2	0.47	2.09	25	8
1:A:8:A:H5''	1:A:9:A:OP1	0.47	2.10	24	1
1:A:6:G:O2'	1:A:7:A:OP1	0.47	2.30	21	1
1:A:2:C:C4	1:A:3:C:C5	0.47	3.03	11	1
2:B:7:ARG:NH2	2:B:11:ARG:CD	0.47	2.78	6	10
1:A:14:G:C2'	1:A:15:C:H5''	0.47	2.40	11	2
1:A:9:A:C8	1:A:9:A:OP1	0.46	2.67	25	8
1:A:6:G:C4	1:A:8:A:OP2	0.46	2.68	7	4
1:A:5:U:OP2	2:B:11:ARG:CZ	0.46	2.62	1	1
1:A:4:C:C5	2:B:7:ARG:HD2	0.46	2.45	23	5
1:A:4:C:C5	2:B:7:ARG:CD	0.46	2.99	11	2
1:A:5:U:C2	1:A:12:G:C2	0.46	3.03	2	2
1:A:6:G:H2'	1:A:8:A:OP2	0.46	2.09	21	1
1:A:7:A:O2'	1:A:8:A:O4'	0.46	2.33	28	1
1:A:3:C:C2'	1:A:4:C:H5'	0.46	2.41	1	1
1:A:4:C:H2'	1:A:5:U:H6	0.46	1.70	22	6
1:A:14:G:O6	1:A:15:C:N4	0.46	2.49	24	2
1:A:9:A:OP1	2:B:8:ARG:HG2	0.45	2.12	24	1
1:A:8:A:O5'	2:B:15:GLN:CG	0.45	2.64	28	1
1:A:8:A:O5'	2:B:15:GLN:CD	0.45	2.54	28	1
2:B:13:GLU:O	2:B:17:GLN:HB2	0.45	2.12	11	1
1:A:1:G:HO2'	1:A:2:C:P	0.45	2.33	13	2
1:A:5:U:OP2	2:B:7:ARG:NH1	0.45	2.50	26	2
1:A:2:C:C2'	1:A:3:C:O5'	0.45	2.65	11	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:5:U:H5''	2:B:7:ARG:NH2	0.44	2.28	5	1
1:A:13:G:O2'	1:A:14:G:C5'	0.44	2.66	3	5
1:A:8:A:O3'	2:B:8:ARG:NH1	0.44	2.51	29	6
1:A:9:A:C6	2:B:15:GLN:OE1	0.44	2.71	24	1
1:A:9:A:OP2	2:B:8:ARG:CD	0.44	2.65	24	1
1:A:13:G:C2	1:A:14:G:C8	0.44	3.05	13	1
2:B:11:ARG:O	2:B:14:LYS:N	0.44	2.51	21	8
1:A:2:C:HO2'	1:A:3:C:P	0.44	2.36	4	1
2:B:15:GLN:O	2:B:19:LYS:N	0.44	2.34	28	1
1:A:7:A:O2'	1:A:8:A:H5'	0.44	2.13	29	1
1:A:4:C:H4'	1:A:5:U:OP1	0.44	2.13	25	2
1:A:12:G:C2	1:A:13:G:C5	0.44	3.05	12	1
1:A:8:A:OP1	2:B:15:GLN:NE2	0.43	2.51	21	1
1:A:6:G:O2'	1:A:8:A:N7	0.43	2.47	15	1
1:A:9:A:OP2	2:B:8:ARG:HD2	0.43	2.13	24	1
1:A:8:A:O2'	1:A:9:A:OP1	0.43	2.36	5	1
1:A:7:A:H5''	2:B:14:LYS:HB3	0.43	1.90	3	4
1:A:8:A:OP1	2:B:11:ARG:CB	0.43	2.67	20	10
1:A:6:G:N2	1:A:10:A:H8	0.43	2.12	15	1
1:A:5:U:C2'	1:A:6:G:C5'	0.43	2.97	16	8
1:A:6:G:C1'	1:A:8:A:N7	0.43	2.82	24	1
1:A:11:A:N6	1:A:12:G:C6	0.43	2.87	3	4
1:A:8:A:OP1	2:B:11:ARG:HB3	0.43	2.13	27	6
1:A:9:A:N6	2:B:19:LYS:HE2	0.43	2.29	1	1
1:A:7:A:O3'	2:B:15:GLN:HG3	0.43	2.14	24	1
1:A:11:A:N6	1:A:12:G:C5	0.43	2.87	20	2
1:A:6:G:O6	2:B:8:ARG:NH2	0.43	2.52	21	1
1:A:9:A:N6	2:B:15:GLN:OE1	0.43	2.52	24	1
1:A:3:C:C4	1:A:4:C:N4	0.42	2.86	29	1
2:B:7:ARG:O	2:B:10:ARG:N	0.42	2.52	23	8
1:A:8:A:N6	1:A:10:A:C5	0.42	2.88	5	3
1:A:6:G:C8	1:A:8:A:OP2	0.42	2.73	12	1
1:A:5:U:O2'	1:A:6:G:C5'	0.42	2.67	14	4
2:B:18:TRP:CE3	2:B:19:LYS:CA	0.42	3.03	27	21
1:A:7:A:O2'	2:B:15:GLN:HG2	0.42	2.15	21	1
1:A:10:A:H2'	1:A:11:A:C5'	0.42	2.45	17	1
2:B:15:GLN:O	2:B:19:LYS:HG3	0.42	2.15	26	2
2:B:7:ARG:NH1	2:B:11:ARG:NE	0.42	2.68	19	1
1:A:8:A:O5'	2:B:8:ARG:HD2	0.42	2.14	25	5
1:A:11:A:N1	1:A:12:G:C5	0.42	2.87	21	1
1:A:5:U:O4	1:A:6:G:O6	0.42	2.38	5	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:9:A:OP2	2:B:8:ARG:CG	0.42	2.68	16	3
1:A:8:A:O3'	1:A:9:A:H3'	0.42	2.15	24	1
1:A:8:A:OP2	2:B:11:ARG:HB3	0.41	2.15	28	1
1:A:2:C:H2'	1:A:3:C:O5'	0.41	2.15	8	4
2:B:13:GLU:CG	2:B:14:LYS:N	0.41	2.83	22	1
2:B:18:TRP:CZ3	2:B:19:LYS:HG2	0.41	2.50	5	1
1:A:8:A:H5''	1:A:9:A:OP2	0.41	2.14	26	1
1:A:5:U:O2'	1:A:6:G:H5'	0.41	2.15	26	2
1:A:10:A:H3'	1:A:11:A:C8	0.41	2.51	4	1
2:B:18:TRP:HE3	2:B:19:LYS:N	0.41	2.14	20	4
1:A:1:G:C6	1:A:2:C:N4	0.41	2.88	24	1
1:A:6:G:O2'	1:A:7:A:C8	0.41	2.73	6	1
1:A:1:G:C6	1:A:2:C:C4	0.40	3.09	19	1
1:A:6:G:O6	2:B:7:ARG:NH2	0.40	2.54	2	1
1:A:6:G:H21	1:A:10:A:H8	0.40	1.58	15	1
2:B:7:ARG:NH1	2:B:11:ARG:CZ	0.40	2.85	14	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	
2	B	17/35 (49%)	16±0 (94±0%)	1±0 (6±0%)	0±0 (0±0%)	100	100
All	All	493/1015 (49%)	464 (94%)	29 (6%)	0 (0%)	100	100

There are no Ramachandran outliers.

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	
2	B	13/28 (46%)	10±1 (79±7%)	3±1 (21±7%)	4	33
All	All	377/812 (46%)	298 (79%)	79 (21%)	4	33

All 6 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)
2	B	15	GLN	29
2	B	10	ARG	19
2	B	8	ARG	12
2	B	7	ARG	9
2	B	17	GLN	7
2	B	13	GLU	3

6.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers	Suiteness
1	A	15/15 (100%)	9±1 (58±6%)	5±1 (31±7%)	0.29±0.07
All	All	423/435 (97%)	254 (60%)	135 (32%)	0.29

The overall RNA backbone suiteness is 0.29.

All unique RNA backbone outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
1	A	9	A	29
1	A	12	G	29
1	A	10	A	29
1	A	6	G	29
1	A	14	G	24
1	A	11	A	24
1	A	4	C	23
1	A	2	C	20
1	A	3	C	19
1	A	5	U	13
1	A	13	G	11
1	A	15	C	2
1	A	8	A	1
1	A	7	A	1

All unique RNA pucker outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
1	A	8	A	28
1	A	3	C	22
1	A	1	G	17
1	A	13	G	16
1	A	11	A	15
1	A	4	C	12
1	A	5	U	11
1	A	12	G	8
1	A	14	G	2
1	A	6	G	1
1	A	10	A	1
1	A	2	C	1
1	A	7	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 carbohydrates 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

No chemical shift data were provided