Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:43 PM GMT

PDB ID	:	1YFG
Title	:	YEAST INITIATOR TRNA
Authors	:	Basavappa, R.; Sigler, P.B.
Deposited on	:	1997-05-08
Resolution	:	3.00 Å(reported)

This is a Full wwPDB X-ray 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/XrayValidationReportHelp
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:

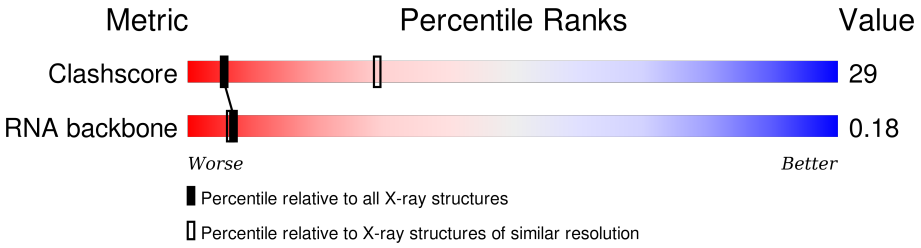
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	1912 (3.00-3.00)
RNA backbone	2183	1036 (3.40-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	75	. 28% 37% 32%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1639 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called YEAST INITIATOR TRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	A	75	1639	734	298	531	76	0	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

- Molecule 1: YEAST INITIATOR TRNA

Chain A: 28% 37% 32%

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α, β, γ	113.72Å 113.72Å 136.54Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 3.00	Depositor
% Data completeness (in resolution range)	81.7 (10.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
Refinement program	NUCLIN, PROFFT (MODIFIED BY Z.OTWINOWSKI)	Depositor
R, R_{free}	0.215 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1639	wwPDB-VP
Average B, all atoms (Å^2)	57.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: H2U, T6A, 2MG, 5MC, 1MA, M2G, 7MG, 1MG, RIA

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 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	1.77	16/1530 (1.0%)	3.19	245/2380 (10.3%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	41	C	O3'-P	-8.59	1.50	1.61
1	A	18	G	O3'-P	7.16	1.69	1.61
1	A	65	G	P-O5'	-6.79	1.52	1.59
1	A	54	A	O3'-P	6.37	1.68	1.61
1	A	21	A	N9-C4	6.29	1.41	1.37
1	A	41	C	C5'-C4'	6.02	1.58	1.51
1	A	18	G	C6-N1	-5.55	1.35	1.39
1	A	57	G	C6-N1	-5.33	1.35	1.39
1	A	53	G	C4'-O4'	-5.29	1.38	1.45
1	A	63	G	P-O5'	5.25	1.65	1.59
1	A	1	A	OP3-P	5.20	1.67	1.61
1	A	71	C	C4'-O4'	-5.20	1.38	1.45
1	A	63	G	C4'-O4'	-5.19	1.38	1.45
1	A	8	U	C4-O4	5.13	1.27	1.23
1	A	45	U	C4-O4	5.12	1.27	1.23
1	A	72	U	C2'-C1'	-5.03	1.47	1.53

All (245) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	37	T6A	P-O3'-C3'	26.06	150.97	119.70
1	A	38	A	P-O3'-C3'	20.33	144.09	119.70
1	A	76	A	O4'-C1'-N9	20.29	124.44	108.20
1	A	29	G	P-O3'-C3'	17.69	140.92	119.70
1	A	19	G	O4'-C1'-N9	-16.88	94.70	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	34	C	P-O3'-C3'	16.25	139.20	119.70
1	A	74	C	P-O3'-C3'	14.99	137.69	119.70
1	A	19	G	O5'-P-OP2	13.91	127.39	110.70
1	A	18	G	O4'-C1'-N9	13.87	119.29	108.20
1	A	63	G	P-O5'-C5'	-13.31	99.60	120.90
1	A	50	U	P-O3'-C3'	-12.76	104.39	119.70
1	A	54	A	P-O3'-C3'	-12.76	104.39	119.70
1	A	21	A	O4'-C1'-N9	12.41	118.13	108.20
1	A	35	A	P-O3'-C3'	12.13	134.25	119.70
1	A	38	A	C6-N1-C2	12.08	125.85	118.60
1	A	15	G	P-O5'-C5'	-12.00	101.70	120.90
1	A	32	C	P-O5'-C5'	11.90	139.94	120.90
1	A	39	C	P-O3'-C3'	11.85	133.92	119.70
1	A	18	G	P-O5'-C5'	-11.30	102.82	120.90
1	A	71	C	O5'-P-OP2	11.25	124.20	110.70
1	A	51	C	O5'-P-OP1	10.99	123.89	110.70
1	A	50	U	C5-C4-O4	-10.74	119.46	125.90
1	A	55	U	P-O5'-C5'	-10.61	103.92	120.90
1	A	59	A	OP1-P-OP2	-10.60	103.70	119.60
1	A	40	C	O5'-P-OP2	10.58	123.40	110.70
1	A	69	C	OP1-P-OP2	-10.56	103.76	119.60
1	A	60	A	OP1-P-OP2	-10.43	103.95	119.60
1	A	45	U	O4'-C1'-N1	10.42	116.53	108.20
1	A	38	A	N1-C2-N3	-10.07	124.26	129.30
1	A	24	G	P-O3'-C3'	10.00	131.70	119.70
1	A	57	G	P-O5'-C5'	-9.84	105.15	120.90
1	A	18	G	P-O3'-C3'	-9.79	107.95	119.70
1	A	69	C	P-O3'-C3'	-9.73	108.03	119.70
1	A	16	H2U	P-O3'-C3'	-9.69	108.08	119.70
1	A	72	U	P-O3'-C3'	-9.61	108.17	119.70
1	A	11	C	P-O3'-C3'	-9.54	108.25	119.70
1	A	50	U	O5'-P-OP2	9.43	122.01	110.70
1	A	52	G	P-O5'-C5'	-9.42	105.83	120.90
1	A	32	C	P-O3'-C3'	9.36	130.93	119.70
1	A	60	A	P-O5'-C5'	-9.20	106.18	120.90
1	A	38	A	C5-C6-N1	-9.06	113.17	117.70
1	A	15	G	O5'-P-OP2	9.03	121.54	110.70
1	A	56	C	P-O5'-C5'	-8.96	106.57	120.90
1	A	32	C	N1-C2-O2	8.93	124.26	118.90
1	A	72	U	C2-N3-C4	-8.80	121.72	127.00
1	A	15	G	N9-C1'-C2'	-8.73	102.39	112.00
1	A	50	U	N3-C4-C5	8.72	119.83	114.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	74	C	O4'-C1'-N1	8.71	115.17	108.20
1	A	36	U	C2-N3-C4	-8.66	121.81	127.00
1	A	62	C	O5'-P-OP1	8.61	121.04	110.70
1	A	55	U	N3-C4-C5	8.59	119.75	114.60
1	A	5	C	P-O3'-C3'	-8.46	109.55	119.70
1	A	72	U	P-O5'-C5'	-8.45	107.38	120.90
1	A	33	U	C2-N3-C4	-8.43	121.94	127.00
1	A	50	U	C2-N3-C4	-8.39	121.97	127.00
1	A	51	C	P-O3'-C3'	-8.39	109.63	119.70
1	A	65	G	OP1-P-OP2	-8.38	107.03	119.60
1	A	36	U	N3-C4-C5	8.38	119.63	114.60
1	A	69	C	O5'-P-OP2	8.36	120.73	110.70
1	A	8	U	C2-N3-C4	-8.35	121.99	127.00
1	A	55	U	O5'-P-OP2	8.31	120.67	110.70
1	A	56	C	OP1-P-OP2	-8.29	107.16	119.60
1	A	12	G	O5'-P-OP1	8.23	120.58	110.70
1	A	31	G	O3'-P-O5'	8.17	119.53	104.00
1	A	52	G	O4'-C1'-N9	-8.15	101.68	108.20
1	A	45	U	C2-N3-C4	-8.12	122.13	127.00
1	A	68	G	OP1-P-OP2	-8.08	107.47	119.60
1	A	72	U	N3-C4-C5	8.04	119.42	114.60
1	A	73	A	N1-C2-N3	-7.94	125.33	129.30
1	A	75	C	N3-C4-C5	-7.94	118.72	121.90
1	A	42	U	N3-C4-C5	7.92	119.35	114.60
1	A	33	U	N3-C4-C5	7.91	119.34	114.60
1	A	6	C	P-O5'-C5'	-7.90	108.27	120.90
1	A	55	U	C2-N3-C4	-7.90	122.26	127.00
1	A	66	C	O4'-C1'-N1	7.88	114.51	108.20
1	A	24	G	N9-C1'-C2'	-7.86	103.36	112.00
1	A	23	C	O5'-P-OP2	7.78	120.04	110.70
1	A	62	C	OP2-P-O3'	7.75	122.25	105.20
1	A	49	5MC	P-O3'-C3'	-7.70	110.47	119.70
1	A	70	G	P-O5'-C5'	-7.68	108.60	120.90
1	A	19	G	OP1-P-OP2	-7.67	108.09	119.60
1	A	45	U	N3-C4-C5	7.64	119.19	114.60
1	A	53	G	P-O5'-C5'	-7.63	108.69	120.90
1	A	68	G	O4'-C1'-N9	7.55	114.24	108.20
1	A	52	G	P-O3'-C3'	-7.51	110.68	119.70
1	A	2	G	O5'-P-OP2	7.51	119.71	110.70
1	A	2	G	P-O5'-C5'	-7.49	108.92	120.90
1	A	57	G	O5'-P-OP2	7.46	119.66	110.70
1	A	60	A	O5'-P-OP2	7.39	119.57	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	14	A	O4'-C1'-N9	-7.35	102.32	108.20
1	A	50	U	P-O5'-C5'	-7.31	109.20	120.90
1	A	23	C	P-O5'-C5'	-7.14	109.48	120.90
1	A	22	G	O5'-P-OP2	7.13	119.26	110.70
1	A	2	G	C5-C6-N1	7.13	115.06	111.50
1	A	42	U	C2-N3-C4	-7.10	122.74	127.00
1	A	8	U	N3-C4-C5	7.09	118.86	114.60
1	A	14	A	C5-C6-N1	-7.09	114.15	117.70
1	A	32	C	OP1-P-OP2	-7.05	109.02	119.60
1	A	76	A	C5-C6-N1	-7.03	114.19	117.70
1	A	3	C	P-O3'-C3'	6.98	128.08	119.70
1	A	6	C	O4'-C1'-N1	6.90	113.72	108.20
1	A	43	G	C5-C6-N1	6.76	114.88	111.50
1	A	31	G	N9-C1'-C2'	-6.71	104.62	112.00
1	A	50	U	OP1-P-OP2	-6.69	109.57	119.60
1	A	52	G	O5'-P-OP2	6.68	118.72	110.70
1	A	22	G	C5-C6-N1	6.59	114.80	111.50
1	A	73	A	O5'-P-OP2	6.59	118.61	110.70
1	A	70	G	O5'-P-OP1	6.57	118.59	110.70
1	A	71	C	OP1-P-OP2	-6.55	109.78	119.60
1	A	39	C	OP1-P-OP2	-6.52	109.82	119.60
1	A	22	G	P-O5'-C5'	-6.51	110.48	120.90
1	A	54	A	C5-C6-N1	-6.49	114.45	117.70
1	A	55	U	P-O3'-C3'	6.49	127.49	119.70
1	A	50	U	C5'-C4'-C3'	-6.48	105.63	116.00
1	A	6	C	N3-C4-C5	-6.47	119.31	121.90
1	A	70	G	OP1-P-OP2	-6.47	109.89	119.60
1	A	1	A	OP1-P-OP2	-6.46	109.92	119.60
1	A	66	C	P-O3'-C3'	6.42	127.40	119.70
1	A	33	U	OP1-P-OP2	-6.41	109.99	119.60
1	A	65	G	O5'-P-OP2	6.37	118.34	110.70
1	A	33	U	N1-C2-O2	-6.34	118.36	122.80
1	A	66	C	P-O5'-C5'	-6.34	110.75	120.90
1	A	51	C	O3'-P-O5'	-6.30	92.04	104.00
1	A	39	C	P-O5'-C5'	-6.29	110.84	120.90
1	A	42	U	OP1-P-OP2	-6.28	110.18	119.60
1	A	14	A	N1-C6-N6	6.17	122.30	118.60
1	A	76	A	P-O5'-C5'	-6.17	111.04	120.90
1	A	51	C	P-O5'-C5'	-6.15	111.06	120.90
1	A	58	1MA	OP2-P-O3'	6.15	118.73	105.20
1	A	7	G	C5-C6-N1	6.14	114.57	111.50
1	A	22	G	O4'-C1'-N9	6.14	113.11	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	12	G	P-O3'-C3'	-6.13	112.34	119.70
1	A	65	G	O5'-P-OP1	6.11	118.03	110.70
1	A	63	G	OP1-P-OP2	-6.09	110.46	119.60
1	A	39	C	N3-C4-C5	-6.07	119.47	121.90
1	A	25	C	N3-C4-C5	-6.07	119.47	121.90
1	A	28	A	N1-C2-N3	-6.05	126.27	129.30
1	A	28	A	C6-N1-C2	6.05	122.23	118.60
1	A	27	C	O5'-P-OP2	6.05	117.95	110.70
1	A	45	U	N1-C2-N3	6.04	118.52	114.90
1	A	66	C	N1-C1'-C2'	-6.03	105.37	112.00
1	A	60	A	N1-C2-N3	-6.01	126.29	129.30
1	A	32	C	N3-C2-O2	-5.99	117.71	121.90
1	A	29	G	N9-C1'-C2'	-5.98	105.42	112.00
1	A	55	U	N1-C1'-C2'	-5.98	105.42	112.00
1	A	38	A	OP1-P-OP2	-5.96	110.67	119.60
1	A	12	G	C5-C6-N1	5.95	114.47	111.50
1	A	12	G	OP1-P-OP2	-5.95	110.68	119.60
1	A	72	U	N1-C2-N3	5.94	118.47	114.90
1	A	31	G	C5-C6-N1	5.94	114.47	111.50
1	A	25	C	O4'-C1'-N1	5.93	112.95	108.20
1	A	35	A	C5-C6-N1	-5.90	114.75	117.70
1	A	39	C	O4'-C1'-N1	5.89	112.91	108.20
1	A	31	G	C6-N1-C2	-5.86	121.58	125.10
1	A	33	U	N1-C2-N3	5.84	118.40	114.90
1	A	59	A	O5'-P-OP1	5.82	117.69	110.70
1	A	55	U	O5'-P-OP1	5.80	117.66	110.70
1	A	68	G	O5'-P-OP1	5.80	117.66	110.70
1	A	18	G	C5-C6-N1	5.78	114.39	111.50
1	A	28	A	C5-C6-N1	-5.75	114.83	117.70
1	A	61	C	C6-N1-C2	5.75	122.60	120.30
1	A	73	A	C6-N1-C2	5.73	122.04	118.60
1	A	75	C	C2-N3-C4	5.73	122.77	119.90
1	A	76	A	C6-N1-C2	5.71	122.02	118.60
1	A	34	C	N3-C4-C5	-5.70	119.62	121.90
1	A	72	U	C5-C4-O4	-5.69	122.48	125.90
1	A	35	A	N1-C2-N3	-5.68	126.46	129.30
1	A	11	C	N1-C2-O2	5.65	122.29	118.90
1	A	14	A	C6-N1-C2	5.65	121.99	118.60
1	A	71	C	P-O5'-C5'	-5.65	111.86	120.90
1	A	29	G	C5-C6-N1	5.62	114.31	111.50
1	A	74	C	C2-N3-C4	5.61	122.70	119.90
1	A	5	C	O5'-P-OP1	5.60	117.42	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	30	G	OP1-P-OP2	-5.58	111.22	119.60
1	A	56	C	P-O3'-C3'	-5.56	113.03	119.70
1	A	70	G	C5-C6-N1	5.55	114.28	111.50
1	A	53	G	O5'-P-OP2	5.55	117.36	110.70
1	A	12	G	C6-N1-C2	-5.55	121.77	125.10
1	A	2	G	C6-N1-C2	-5.55	121.77	125.10
1	A	65	G	O5'-C5'-C4'	-5.54	101.17	111.70
1	A	28	A	O4'-C1'-N9	5.54	112.63	108.20
1	A	53	G	C5-C6-N1	5.51	114.25	111.50
1	A	30	G	C5-C6-N1	5.50	114.25	111.50
1	A	59	A	C5-C6-N1	-5.49	114.95	117.70
1	A	35	A	OP1-P-OP2	-5.48	111.37	119.60
1	A	59	A	N1-C6-N6	5.48	121.89	118.60
1	A	52	G	C5-C6-N1	5.45	114.23	111.50
1	A	5	C	N3-C4-C5	-5.44	119.72	121.90
1	A	18	G	N9-C1'-C2'	-5.39	106.07	112.00
1	A	61	C	P-O3'-C3'	-5.39	113.23	119.70
1	A	19	G	C8-N9-C4	-5.38	104.25	106.40
1	A	36	U	OP1-P-OP2	-5.37	111.54	119.60
1	A	67	G	O3'-P-O5'	5.37	114.20	104.00
1	A	51	C	OP1-P-OP2	-5.36	111.57	119.60
1	A	56	C	N3-C4-C5	-5.35	119.76	121.90
1	A	8	U	N1-C2-N3	5.34	118.11	114.90
1	A	7	G	C6-N1-C2	-5.34	121.89	125.10
1	A	22	G	P-O3'-C3'	-5.34	113.29	119.70
1	A	53	G	O4'-C1'-N9	-5.34	103.93	108.20
1	A	53	G	OP1-P-OP2	-5.33	111.61	119.60
1	A	71	C	N1-C1'-C2'	-5.33	106.14	112.00
1	A	43	G	O5'-P-OP2	5.32	117.08	110.70
1	A	66	C	OP2-P-O3'	5.31	116.87	105.20
1	A	56	C	OP1-P-O3'	-5.29	93.55	105.20
1	A	43	G	C6-N1-C2	-5.29	121.92	125.10
1	A	54	A	C6-N1-C2	5.28	121.77	118.60
1	A	3	C	N1-C1'-C2'	-5.28	106.19	112.00
1	A	4	G	C5-C6-N1	5.27	114.14	111.50
1	A	55	U	C5-C4-O4	-5.27	122.74	125.90
1	A	57	G	C5-C6-O6	-5.27	125.44	128.60
1	A	70	G	C5-C6-O6	-5.25	125.45	128.60
1	A	53	G	C6-N1-C2	-5.25	121.95	125.10
1	A	41	C	O4'-C1'-N1	5.24	112.39	108.20
1	A	73	A	C5-C6-N1	-5.23	115.09	117.70
1	A	1	A	N1-C2-N3	-5.22	126.69	129.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	4	G	O5'-P-OP1	5.21	116.96	110.70
1	A	36	U	O5'-P-OP2	5.21	116.95	110.70
1	A	52	G	OP1-P-OP2	-5.20	111.81	119.60
1	A	54	A	O5'-P-OP1	5.20	116.94	110.70
1	A	11	C	N3-C4-C5	-5.17	119.83	121.90
1	A	27	C	N3-C4-C5	-5.15	119.84	121.90
1	A	66	C	N3-C4-C5	-5.14	119.84	121.90
1	A	57	G	C5-C6-N1	5.14	114.07	111.50
1	A	30	G	C6-N1-C2	-5.13	122.02	125.10
1	A	16	H2U	OP1-P-O3'	5.13	116.48	105.20
1	A	30	G	O4'-C1'-N9	-5.12	104.10	108.20
1	A	5	C	OP1-P-OP2	-5.12	111.93	119.60
1	A	12	G	P-O5'-C5'	-5.12	112.72	120.90
1	A	52	G	C6-N1-C2	-5.11	122.03	125.10
1	A	51	C	OP1-P-O3'	5.10	116.42	105.20
1	A	36	U	N1-C2-N3	5.09	117.95	114.90
1	A	44	A	N1-C6-N6	5.09	121.65	118.60
1	A	42	U	O5'-P-OP2	5.09	116.81	110.70
1	A	30	G	C5-C6-O6	-5.08	125.55	128.60
1	A	44	A	OP1-P-OP2	-5.07	111.99	119.60
1	A	65	G	C6-N1-C2	-5.07	122.06	125.10
1	A	22	G	C6-N1-C2	-5.06	122.06	125.10
1	A	31	G	O4'-C1'-N9	5.05	112.24	108.20
1	A	18	G	O4'-C1'-C2'	-5.05	100.75	105.80
1	A	76	A	N9-C1'-C2'	-5.02	106.48	112.00
1	A	67	G	OP1-P-OP2	-5.01	112.08	119.60
1	A	42	U	O4'-C1'-N1	5.01	112.20	108.20
1	A	72	U	O5'-P-OP2	5.01	116.71	110.70
1	A	7	G	C5-C6-O6	-5.00	125.60	128.60
1	A	65	G	OP2-P-O3'	5.00	116.21	105.20

There are no chirality outliers.

There are no planarity outliers.

5.2 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 the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1639	0	848	72	0
All	All	1639	0	848	72	0

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 29.

All (72) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:C:H2'	1:A:42:U:C6	2.07	0.90
1:A:47:H2U:H2'	1:A:47:H2U:O2	1.75	0.86
1:A:50:U:H5'	1:A:51:C:OP2	1.77	0.83
1:A:59:A:O2'	1:A:60:A:H5'	1.84	0.78
1:A:38:A:H2'	1:A:39:C:H6	1.48	0.77
1:A:32:C:H2'	1:A:32:C:O2	1.85	0.75
1:A:49:5MC:H2'	1:A:49:5MC:O2	1.87	0.74
1:A:38:A:O2'	1:A:39:C:H5'	1.87	0.74
1:A:35:A:H2'	1:A:36:U:O4'	1.91	0.70
1:A:7:G:C5	1:A:49:5MC:HM52	2.27	0.69
1:A:38:A:H2'	1:A:39:C:C6	2.28	0.67
1:A:38:A:H5'	1:A:38:A:H8	1.60	0.67
1:A:39:C:N3	1:A:40:C:C5	2.64	0.65
1:A:35:A:H2'	1:A:36:U:C6	2.32	0.65
1:A:43:G:H2'	1:A:43:G:N3	2.13	0.63
1:A:38:A:C2	1:A:39:C:C2	2.88	0.62
1:A:10:2MG:H2'	1:A:10:2MG:N3	2.16	0.61
1:A:14:A:C2'	1:A:15:G:H5'	2.31	0.60
1:A:38:A:C8	1:A:38:A:H5'	2.37	0.59
1:A:35:A:C2	1:A:36:U:C2	2.90	0.59
1:A:11:C:O2	1:A:11:C:H2'	2.02	0.59
1:A:54:A:H5"	1:A:55:U:OP2	2.02	0.59
1:A:35:A:H2'	1:A:36:U:H6	1.68	0.58
1:A:23:C:C2'	1:A:24:G:H5'	2.34	0.57
1:A:13:C:H2'	1:A:14:A:H5"	1.86	0.57
1:A:56:C:O2	1:A:56:C:H2'	2.04	0.57
1:A:24:G:O2'	1:A:25:C:H5'	2.05	0.56
1:A:38:A:N3	1:A:39:C:C6	2.74	0.56
1:A:44:A:OP2	1:A:44:A:H8	1.91	0.53
1:A:38:A:C4	1:A:39:C:C5	2.96	0.53
1:A:23:C:O2'	1:A:24:G:H5'	2.08	0.53
1:A:49:5MC:C2'	1:A:49:5MC:O2	2.57	0.53
1:A:61:C:H2'	1:A:61:C:O2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:U:H2'	1:A:50:U:O2	2.08	0.52
1:A:56:C:H5"	1:A:57:G:OP2	2.10	0.52
1:A:40:C:C2'	1:A:41:C:H5'	2.40	0.51
1:A:59:A:HO2'	1:A:60:A:H5'	1.75	0.51
1:A:75:C:H2'	1:A:76:A:O4'	2.11	0.51
1:A:38:A:C4	1:A:39:C:C6	2.99	0.51
1:A:39:C:C2	1:A:40:C:C5	2.98	0.51
1:A:35:A:C6	1:A:36:U:C4	3.00	0.50
1:A:14:A:H2'	1:A:15:G:H5'	1.95	0.49
1:A:40:C:H2'	1:A:41:C:O4'	2.13	0.49
1:A:33:U:O2	1:A:35:A:N7	2.46	0.48
1:A:31:G:C2	1:A:40:C:C4	3.02	0.47
1:A:12:G:OP2	1:A:12:G:H8	1.97	0.47
1:A:13:C:C2'	1:A:14:A:H5"	2.43	0.47
1:A:39:C:C2	1:A:40:C:C6	3.03	0.47
1:A:24:G:H3'	1:A:24:G:C8	2.51	0.46
1:A:31:G:N1	1:A:40:C:C4	2.84	0.46
1:A:25:C:N4	1:A:26:M2G:C5	2.85	0.45
1:A:31:G:H2'	1:A:32:C:H6	1.82	0.44
1:A:31:G:H2'	1:A:32:C:C6	2.52	0.44
1:A:9:1MG:HN21	1:A:9:1MG:HM11	1.52	0.44
1:A:38:A:H2'	1:A:39:C:O4'	2.17	0.44
1:A:26:M2G:HM22	1:A:44:A:H2	1.82	0.43
1:A:58:1MA:H1'	1:A:60:A:N7	2.34	0.43
1:A:24:G:C2'	1:A:25:C:H5'	2.49	0.43
1:A:31:G:H1	1:A:39:C:H42	1.65	0.43
1:A:15:G:C2	1:A:59:A:C2	3.07	0.43
1:A:5:C:H2'	1:A:6:C:O4'	2.19	0.43
1:A:61:C:C2'	1:A:61:C:O2	2.64	0.42
1:A:11:C:C2'	1:A:11:C:O2	2.67	0.42
1:A:25:C:C4	1:A:26:M2G:C8	3.08	0.42
1:A:35:A:C5	1:A:36:U:C5	3.08	0.42
1:A:35:A:C4	1:A:36:U:C6	3.07	0.42
1:A:24:G:C8	1:A:24:G:C3'	3.03	0.41
1:A:27:C:O2'	1:A:28:A:H5'	2.19	0.41
1:A:56:C:O2	1:A:56:C:C2'	2.66	0.41
1:A:4:G:H2'	1:A:5:C:C6	2.56	0.41
1:A:71:C:O2	1:A:71:C:H2'	2.21	0.41
1:A:34:C:OP1	1:A:34:C:H6	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

There are no protein molecules in this entry.

5.3.2 Protein sidechains

There are no protein molecules in this entry.

5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	73/75 (97%)	37 (50%)	7 (9%)

All (37) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	5	C
1	A	9	1MG
1	A	11	C
1	A	12	G
1	A	13	C
1	A	14	A
1	A	16	H2U
1	A	18	G
1	A	19	G
1	A	21	A
1	A	22	G
1	A	24	G
1	A	29	G
1	A	30	G
1	A	32	C
1	A	33	U
1	A	34	C
1	A	35	A
1	A	36	U
1	A	37	T6A
1	A	38	A
1	A	39	C
1	A	40	C
1	A	41	C
1	A	42	U

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Mol	Chain	Res	Type
1	A	43	G
1	A	44	A
1	A	47	H2U
1	A	48	5MC
1	A	49	5MC
1	A	50	U
1	A	54	A
1	A	56	C
1	A	58	1MA
1	A	61	C
1	A	64	RIA
1	A	75	C

All (7) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	35	A
1	A	36	U
1	A	37	T6A
1	A	38	A
1	A	39	C
1	A	47	H2U
1	A	74	C

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

11 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	2MG	A	10	1	17,26,27	1.18	2 (11%)	21,38,41	3.51	6 (28%)
1	H2U	A	16	1	17,21,22	0.98	1 (5%)	23,30,33	1.18	2 (8%)
1	M2G	A	26	1	17,27,28	1.35	2 (11%)	22,40,43	2.90	7 (31%)
1	T6A	A	37	1	22,34,35	1.81	4 (18%)	25,49,52	3.89	9 (36%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
1	7MG	A	46	1	19,26,27	1.81	2 (10%)	24,39,42	1.96	5 (20%)
1	H2U	A	47	1	17,21,22	1.09	0	23,30,33	2.04	6 (26%)
1	5MC	A	48	1	13,22,23	1.28	3 (23%)	15,32,35	1.02	0
1	5MC	A	49	1	13,22,23	1.25	1 (7%)	15,32,35	1.44	2 (13%)
1	1MA	A	58	1	14,25,26	2.40	3 (21%)	15,37,40	2.56	9 (60%)
1	RIA	A	64	1	29,38,39	1.25	3 (10%)	38,57,60	2.07	9 (23%)
1	1MG	A	9	1	16,26,27	0.81	0	19,39,42	1.89	4 (21%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	2MG	A	10	1	-	0/5/27/28	0/3/3/3
1	H2U	A	16	1	-	0/7/38/39	0/2/2/2
1	M2G	A	26	1	-	0/7/29/30	0/3/3/3
1	T6A	A	37	1	-	0/15/41/42	0/3/3/3
1	7MG	A	46	1	-	0/7/37/38	0/3/3/3
1	H2U	A	47	1	-	0/7/38/39	0/2/2/2
1	5MC	A	48	1	-	0/3/25/26	0/2/2/2
1	5MC	A	49	1	-	0/3/25/26	0/2/2/2
1	1MA	A	58	1	-	0/3/25/26	0/3/3/3
1	RIA	A	64	1	-	0/13/51/52	0/4/4/4
1	1MG	A	9	1	-	0/3/25/26	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	46	7MG	C8-N9	-6.28	1.36	1.45
1	A	37	T6A	C15-C14	-4.16	1.41	1.51
1	A	37	T6A	O14-C14	-4.07	1.34	1.43
1	A	64	RIA	P'-O3X	-3.89	1.40	1.54
1	A	37	T6A	C12-N11	-2.88	1.40	1.46
1	A	58	1MA	O4'-C4'	-2.88	1.38	1.45
1	A	46	7MG	C8-N7	-2.60	1.31	1.43
1	A	48	5MC	O4'-C4'	-2.27	1.39	1.45
1	A	48	5MC	C6-C5	-2.05	1.34	1.40
1	A	16	H2U	C6-N1	-2.00	1.44	1.47
1	A	64	RIA	C2-N1	2.06	1.37	1.33
1	A	64	RIA	O4'-C1A	2.21	1.44	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	10	2MG	C6-N1	2.27	1.37	1.33
1	A	37	T6A	O4'-C1'	2.34	1.44	1.41
1	A	26	M2G	C6-N1	2.76	1.38	1.33
1	A	26	M2G	O4'-C1'	3.01	1.45	1.41
1	A	48	5MC	O4'-C1'	3.09	1.45	1.41
1	A	10	2MG	O4'-C1'	3.16	1.45	1.41
1	A	49	5MC	O4'-C1'	3.47	1.45	1.41
1	A	58	1MA	O2'-C2'	3.96	1.52	1.43
1	A	58	1MA	O4'-C1'	6.37	1.49	1.41

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	10	2MG	CM2-N2-C2	-12.84	108.58	123.07
1	A	26	M2G	CM1-N2-C2	-8.23	112.88	121.34
1	A	37	T6A	O14-C14-C12	-7.74	92.45	109.08
1	A	26	M2G	C5-C6-N1	-6.91	114.14	123.59
1	A	64	RIA	O2A-C1'-C2'	-6.54	94.15	107.75
1	A	10	2MG	C5-C6-N1	-6.25	115.05	123.59
1	A	46	7MG	C5-C6-N1	-5.63	114.81	123.46
1	A	37	T6A	O10-C10-N6	-5.32	115.50	123.58
1	A	58	1MA	O4'-C1'-N9	-4.93	97.77	108.10
1	A	47	H2U	O2-C2-N3	-4.51	112.65	121.48
1	A	58	1MA	O2'-C2'-C3'	-4.36	97.65	111.83
1	A	26	M2G	C2'-C1'-N9	-3.78	108.52	114.29
1	A	9	1MG	C6-C5-C4	-3.52	117.41	119.93
1	A	26	M2G	CM2-N2-C2	-3.37	117.88	121.34
1	A	10	2MG	C2-N3-C4	-3.15	111.29	115.09
1	A	49	5MC	O2'-C2'-C3'	-3.12	101.67	111.83
1	A	10	2MG	C2'-C1'-N9	-3.00	109.71	114.29
1	A	16	H2U	O2-C2-N3	-3.00	115.61	121.48
1	A	9	1MG	C5-C6-N1	-2.94	114.60	118.33
1	A	26	M2G	C2-N3-C4	-2.77	111.75	115.09
1	A	64	RIA	O2X-P'-O5'	-2.76	98.63	106.56
1	A	10	2MG	C6-C5-C4	-2.75	117.61	120.90
1	A	16	H2U	O2'-C2'-C3'	-2.67	103.14	111.83
1	A	47	H2U	C6-N1-C2	-2.54	118.40	122.23
1	A	46	7MG	N2-C2-N1	-2.53	113.01	117.20
1	A	58	1MA	C2-N3-C4	-2.46	112.59	116.40
1	A	64	RIA	N3-C2-N1	-2.38	127.07	128.89
1	A	9	1MG	N2-C2-N1	-2.29	115.27	118.46
1	A	26	M2G	N3-C2-N2	-2.27	114.58	117.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	64	RIA	O3X-P'-O1X	-2.26	103.30	110.58
1	A	58	1MA	C2'-C1'-N9	-2.17	110.97	114.29
1	A	47	H2U	O2'-C2'-C1'	2.07	117.17	109.94
1	A	58	1MA	C4'-O4'-C1'	2.20	112.13	109.72
1	A	64	RIA	C4-C5-N7	2.23	111.53	109.48
1	A	58	1MA	O4'-C4'-C5'	2.34	117.70	109.32
1	A	37	T6A	O4'-C1'-N9	2.40	113.12	108.10
1	A	58	1MA	O3'-C3'-C4'	2.45	118.42	111.05
1	A	58	1MA	C1'-N9-C4	2.53	130.76	126.94
1	A	49	5MC	N4-C4-N3	2.56	120.66	116.95
1	A	64	RIA	O3X-P'-O5'	2.62	114.10	106.56
1	A	64	RIA	O1'-C1'-C2'	2.65	108.50	104.78
1	A	46	7MG	N2-C2-N3	2.73	121.72	117.20
1	A	47	H2U	O3'-C3'-C2'	2.78	120.87	111.83
1	A	37	T6A	C2'-C1'-N9	3.19	119.17	114.29
1	A	58	1MA	O3'-C3'-C2'	3.29	122.54	111.83
1	A	47	H2U	O3'-C3'-C4'	3.46	121.42	111.05
1	A	37	T6A	N6-C6-N1	3.64	122.29	118.85
1	A	46	7MG	CM7-N7-C8	3.82	131.36	120.52
1	A	26	M2G	CM2-N2-CM1	4.02	129.24	115.96
1	A	37	T6A	C13-C12-N11	4.40	122.32	113.51
1	A	10	2MG	C6-N1-C2	4.54	121.92	115.31
1	A	64	RIA	C1'-O2A-C2A	4.62	130.08	118.01
1	A	46	7MG	C6-N1-C2	4.66	122.41	115.94
1	A	9	1MG	O4'-C1'-N9	5.26	119.12	108.10
1	A	47	H2U	O2-C2-N1	5.63	130.66	123.30
1	A	64	RIA	O5'-P'-O1X	6.02	122.47	107.14
1	A	37	T6A	C6-N6-C10	6.08	138.53	130.39
1	A	37	T6A	N6-C10-N11	6.94	124.86	113.81
1	A	37	T6A	O14-C14-C15	11.70	143.46	109.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	10	2MG	1	0
1	A	26	M2G	3	0
1	A	47	H2U	1	0
1	A	49	5MC	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	58	1MA	1	0
1	A	9	1MG	1	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.