

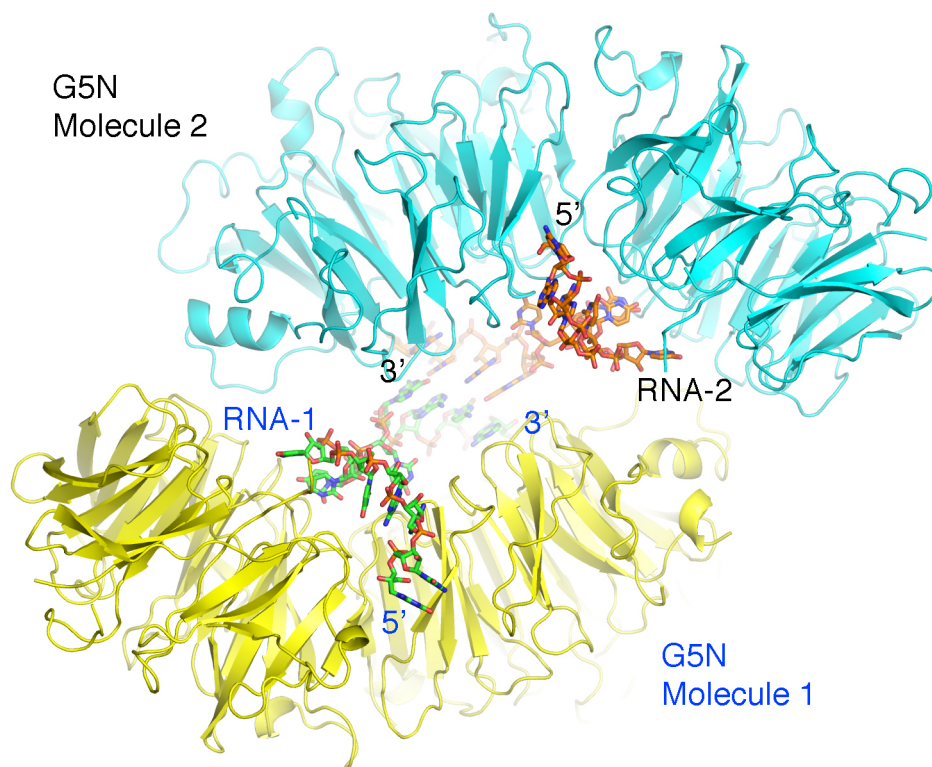
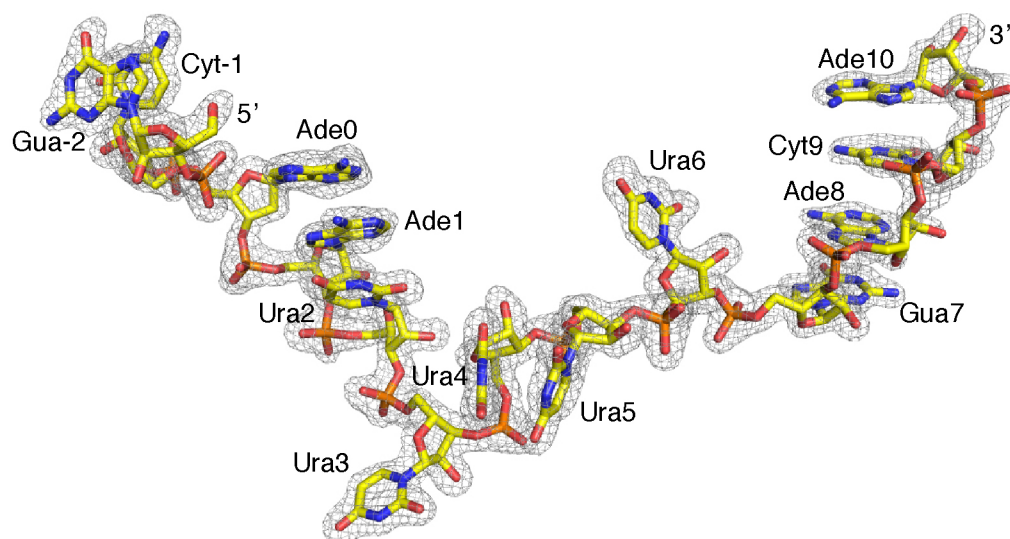
*Supplementary Data*

**Structural basis for snRNA recognition by the double WD40-repeat domain of Gemin5**

Wenxing Jin<sup>1+</sup>, Yi Wang<sup>1,2+</sup>, Chao-Pei Liu<sup>1+</sup>, Na Yang<sup>1,2</sup>, Mingliang Jin<sup>2,3</sup>, Yao Cong<sup>3</sup>,  
Mingzhu Wang<sup>1\*</sup>, Rui-Ming Xu<sup>1,2\*</sup>

<sup>1</sup>National Laboratory of Biomacromolecules, CAS Center for Excellence in Biomacromolecules, Institute of Biophysics, Chinese Academy of Sciences, Beijing 100101, China; <sup>2</sup>University of Chinese Academy of Sciences, Beijing 100049, China; <sup>3</sup>National Center for Protein Science Shanghai, State Key Laboratory of Molecular Biology, Institute of Biochemistry and Cell Biology, Shanghai Institutes for Biological Sciences, Chinese Academy of Sciences, Shanghai 201210, China

Figure S1. Two G5N-RNA complexes in one crystallographic asymmetric unit.

Figure S2. Simulated annealing omit difference electron density map of the bound RNA molecule. The  $F_o - F_c$  map is contoured at  $3\sigma$  level. A stick model of the RNA molecule is superimposed.

Table\_S1. Statistics of crystallographic analyses

	K <sub>2</sub> Pt(NO <sub>2</sub> ) <sub>4</sub> derivative	Native	G5N-13nt RNA (cocrystallization)	G5N-7nt RNA (soak)	G5N-m7GTP (soak)
<b>Data collection</b>					
Space group	P2 <sub>1</sub>	P2 <sub>1</sub>	P2 <sub>1</sub> 2 <sub>1</sub> 2 <sub>1</sub>	P2 <sub>1</sub>	P2 <sub>1</sub>
Cell dimensions					
a, b, c (Å)	59.61, 124.70, 61.64	60.26, 123.59, 61.73	98.65, 107.19, 151.48	59.54, 124.79, 61.02	60.53, 120.91, 61.26
α, β, γ (°)	90, 118.70, 90	90, 119.27, 90	90, 90, 90	90, 117.97, 90	90, 117.42, 90
Wavelength (Å)	0.9786	0.9786	0.9792	0.9785	0.9788
Resolution (Å)	50.00-3.00 (3.11-3.00)	50.00-2.00 (2.07-2.00)	50.00-1.90 (1.97-1.90)	50.00-2.10 (2.18-2.10)	50.00-2.50 (2.59-2.50)
Rmerge	0.143 (0.699)	0.116 (0.504)	0.116 (0.683)	0.090 (0.533)	0.070 (0.609)
I / σI	12 (3.0)	11.5 (2.6)	17.4 (2.8)	13.0 (2.8)	15.9 (2.1)
Completeness (%)	100.0 (100.0)	99.9 (99.9)	99.7 (99.0)	99.7 (99.7)	99.1(99.6)
Redundancy	6.9 (6.8)	4.8 (4.8)	7.9 (7.9)	3.4 (3.4)	3.3(3.3)
Total/Unique reflections	215832/31097	255144/53257	1002669/126392	152592/45763	88710/26965
<b>Refinement</b>					
Resolution (Å)		50.00-2.00 (2.03-2.00)	50.00-1.90 (1.95-1.90)	50.00-2.10 (2.15-2.10)	50.00-2.50 (2.58-2.50)
No. reflections		52884 (2466)	125759 (8413)	45665 (2849)	26936 (2527)
Rwork / Rfree		0.157(0.200)/0.199(0.242)	0.170(0.242)/0.206(0.298)	0.195(0.333)/0.241(0.354)	0.183(0.241)/0.234(0.296)
No. atoms, B-factors (Å <sup>2</sup> )					
Protein		5509, 28.1 Å <sup>2</sup>	10899, 24.0 Å <sup>2</sup>	5378, 34.5 Å <sup>2</sup>	5319, 70.9 Å <sup>2</sup>
RNA			522, 22.7 Å <sup>2</sup>	142, 54.5 Å <sup>2</sup>	
Ligand				6, 29.4 Å <sup>2</sup>	29, 86.8 Å <sup>2</sup>
Water		623, 35.3 Å <sup>2</sup>	1330, 33.3 Å <sup>2</sup>	417, 38.7 Å <sup>2</sup>	33, 58.4 Å <sup>2</sup>
R.m.s deviations					
Bond length (Å)		0.005	0.009	0.005	0.003
Bond angles (°)		0.911	1.256	0.936	0.661
Ramachandran plot					
Favored		96.5%	96.3%	95.5%	97.3%
Allowed		3.5%	3.7%	4.5%	2.6%
Outliers		0	0	0	0.1%