

SUPPLEMENTARY MATERIAL

Structural determinants of the SINE B2 element embedded in the long non-coding RNA activator of translation AS *Uchl1*

Peter Podbevšek^{1,2,3}, Francesca Fasolo¹, Carlotta Bon¹, Laura Cimatti¹, Sabine Reißer⁴, Piero Carninci⁵, Giovanni Bussi⁴, Silvia Zucchelli^{1,6}, Janez Plavec^{2,3,7,*} and Stefano Gustincich^{1,8*}

¹ Area of Neuroscience, SISSA, Trieste, Italy.

² Slovenian NMR Centre, National Institute of Chemistry, Ljubljana, Slovenia.

³ EN-FIST Centre of Excellence, Ljubljana, Slovenia.

⁴ Molecular and Statistical Biophysics, SISSA, Trieste, Italy.

⁵ Division of Genomic Technologies, RIKEN Center for Life Science Technologies, Yokohama, Japan.

⁶ Department of Health Sciences, Università del Piemonte Orientale, Novara, Italy.

⁷ Faculty of Chemistry and Chemical Technology, University of Ljubljana, Ljubljana, Slovenia.

⁸ Department of Neuroscience and Brain Technologies, Italian Institute of Technology, Genova, Italy.

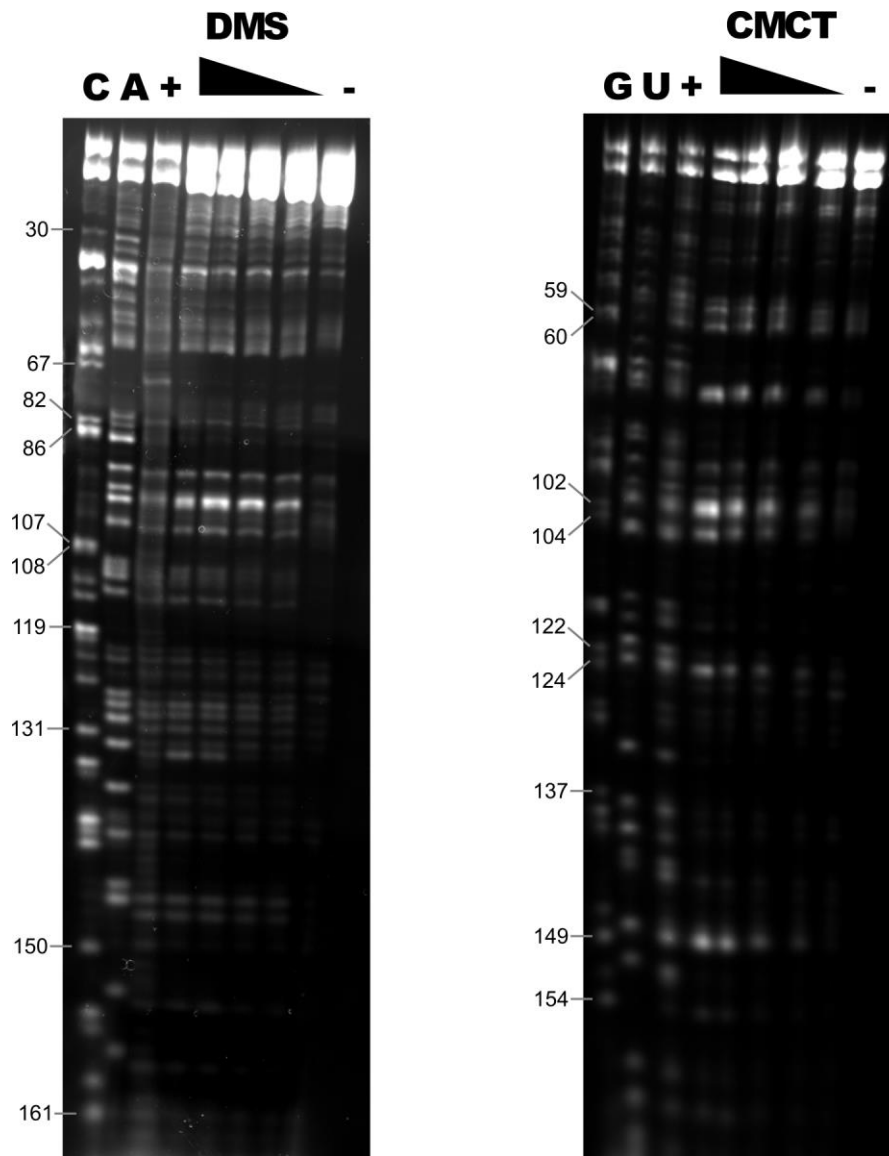


Figure S1 Chemical footprinting of the invSINEB2/183 RNA construct. (left) A fluorescent gel with C and A sequencing lanes with nucleotide numbering. DMS lanes with invSINEB2/183 RNA treated with increasing DMS concentrations with positive (+) and negative (-) controls. (right) A fluorescent gel with G and U sequencing lanes with nucleotide numbering. CMCT lanes with invSINEB2/183 RNA treated with increasing CMCT concentrations with positive (+) and negative (-) controls.

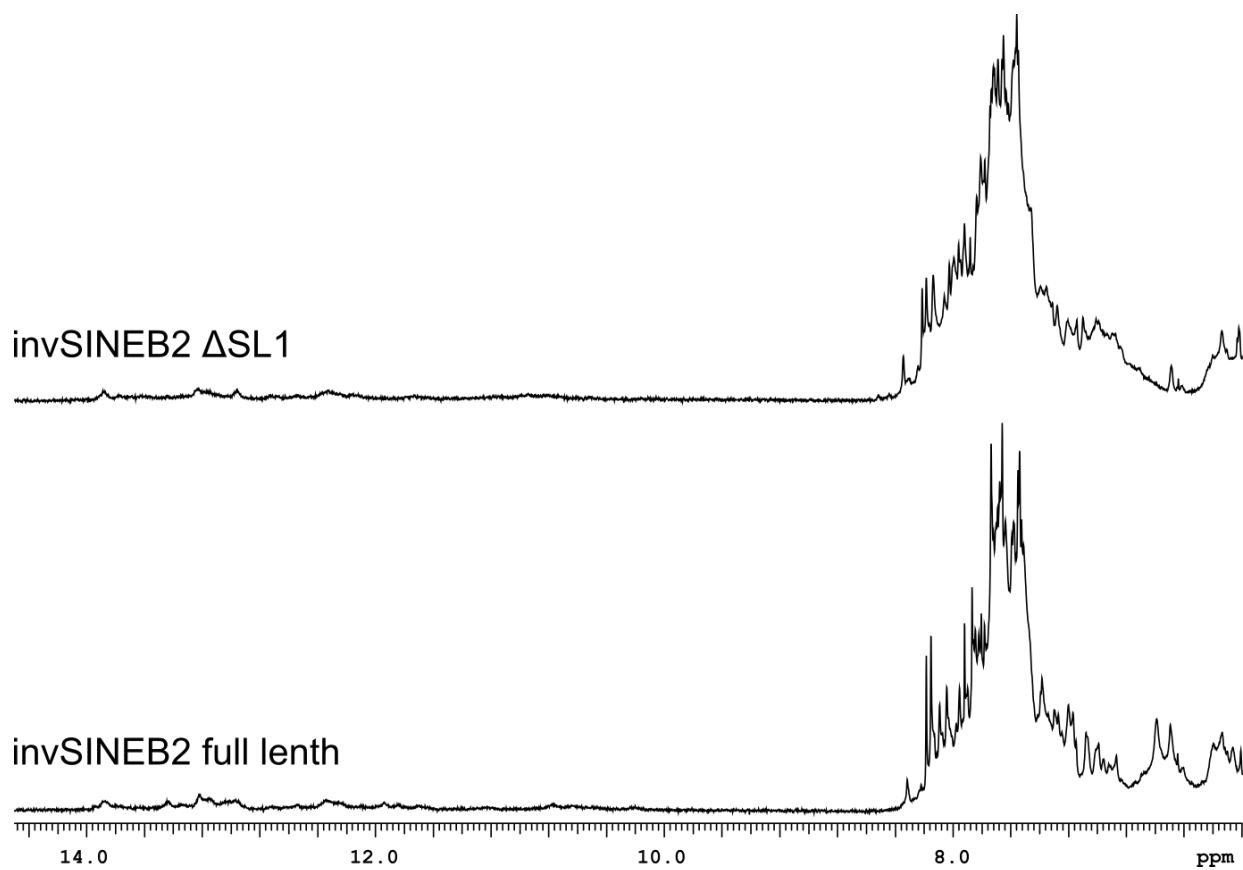


Figure S2 Imino and aromatic regions of ^1H NMR spectra of full length invSINEB2 RNA (183 nt) and the Δ SL1 construct (173 nt) in 5% $2\text{H}_2\text{O}/95\%$ H_2O acquired at 25 $^\circ\text{C}$.

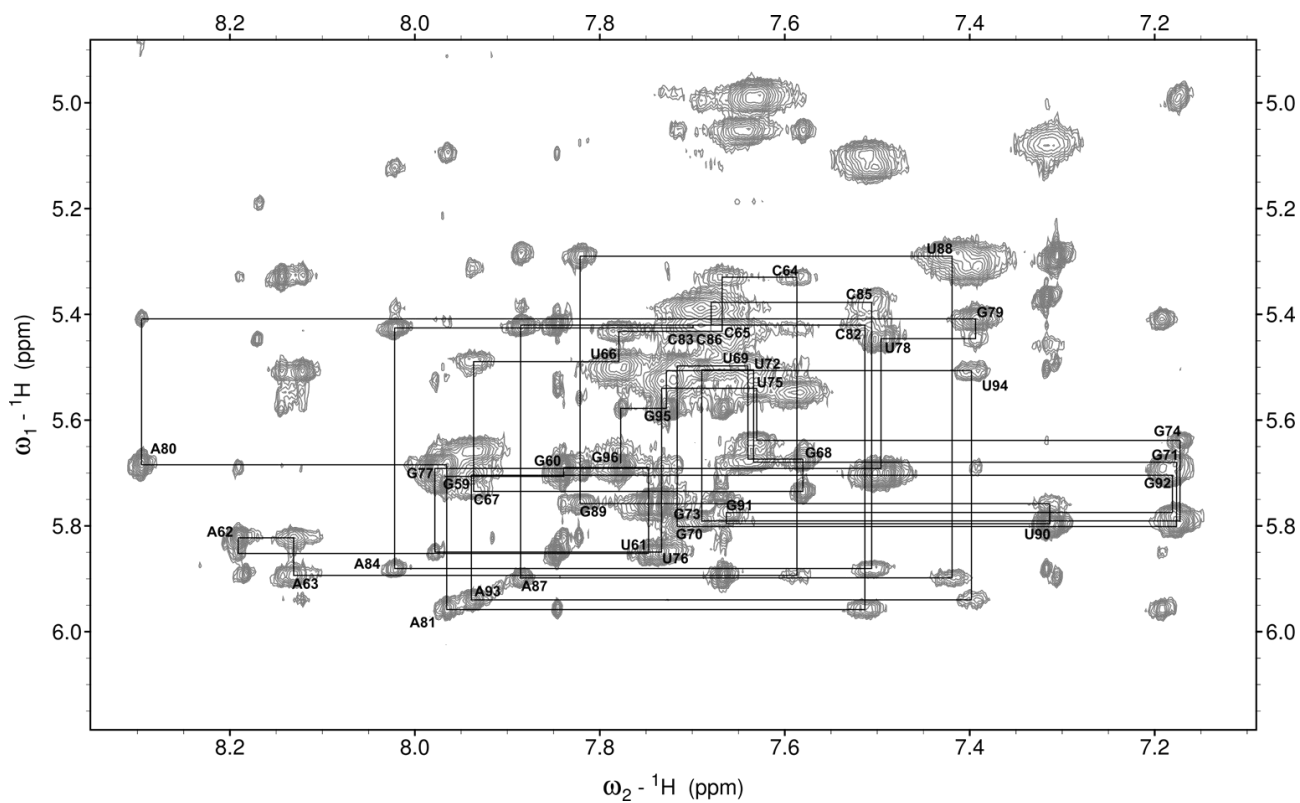


Figure S3 Assignment of relevant resonances. Aromatic-anomeric region of a 2D NOESY NMR spectrum ($\tau_m=250$ ms) of invSINEB2/38. The sequential walk is depicted as a black line.

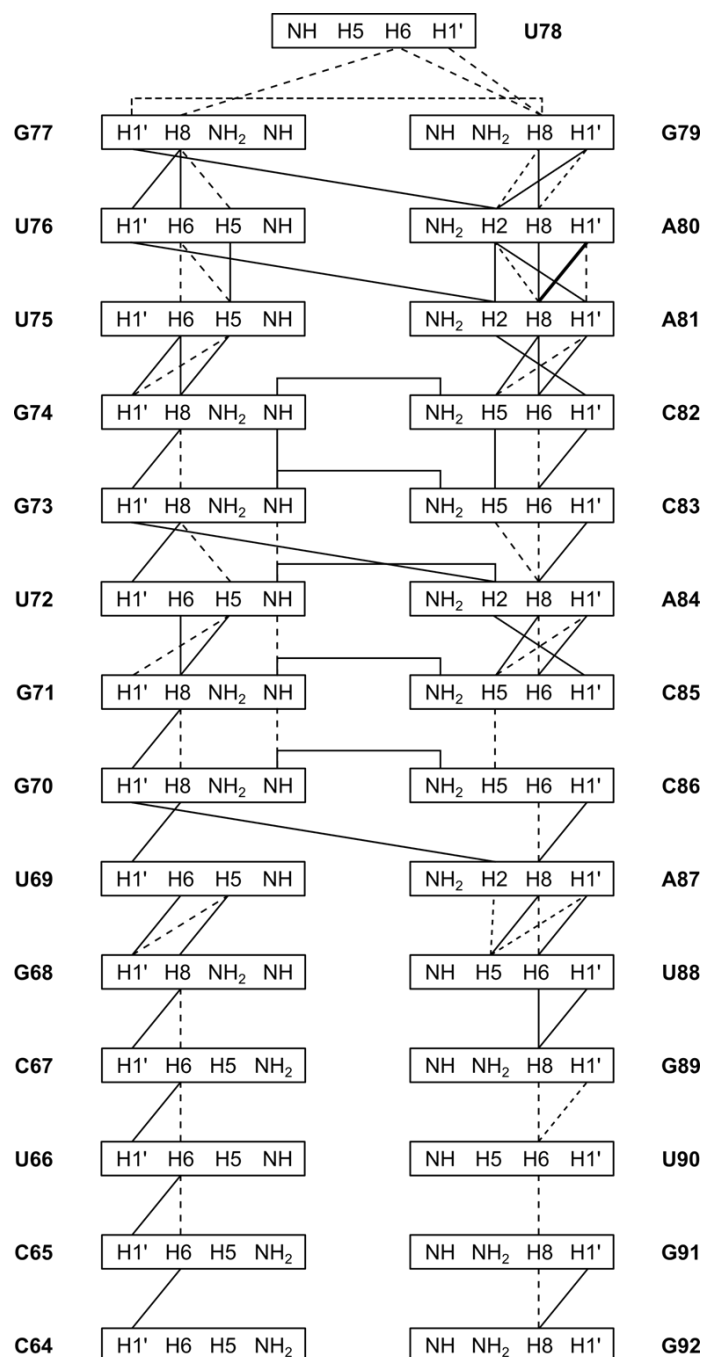


Figure S4 A diagram of strong (bold line) medium (solid line) and weak (dashed line) NOE connectivities for the invSINEB2/38 molecule.

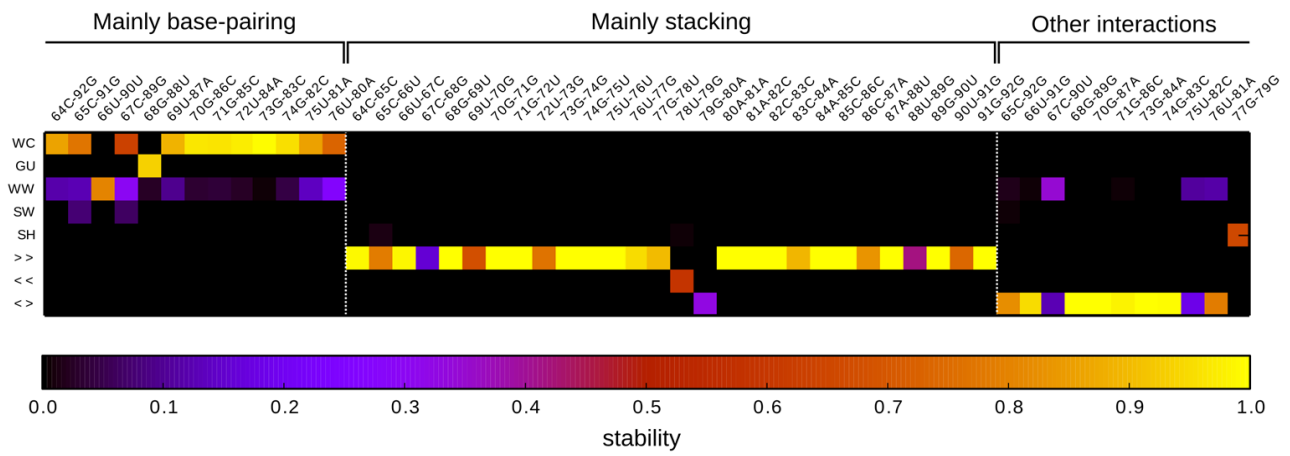


Figure S5 Stability of annotations per pair of bases, calculated with baRNAbA (53) from the MD simulation. WC, GU and WW are interactions between the Watson-Crick edges of both bases. SW is a sugar-edge Watson-Crick-edge interaction, SH is a sugar-edge Hogsteen-edge interaction. >>, << and <> are stackings, in particular >> is the upward stacking found in a canonical double strand, << the downward stacking, and in <> the outward stacking.