Molecular Origami of SCN^-

given information

ElementNames	[(C) (N) (S)]	
dotted	F	
distance	111.198	C^1-N^1
distance	166.468	C^1-S^1
angle	178.101	S^1-C^1-N^1
	277.6	S^1-N^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

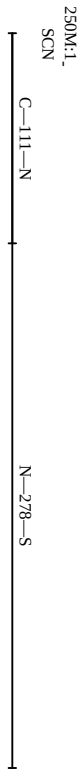
Molecular Origami of SCN^-

!C1
N1
S1
SCN^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.09 offsety 0.9

View -1



Current: (centerx 4.21) (centery 5.90) (scale 250)
%%BoundingBox: 293 139 313 435 actual: 303 149 303 425 center: 303 287 actual size: 0 276
Better: (centerx 4.25) (centery 7.41) (scale 250)
%%BoundingBox: 290 313 310 609 actual: 300 323 300 599 center: 300 461 actual size: 0 276