Molecular Origami of SCN^-

given information

ElementNames	[(C) (N) (S)]	
dotted	F	
distance	115.923	C^1-N^1
distance	161.050	C^1-S^1
angle	179.411	S^1-C^1-N^1
	277.	S^1-N^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

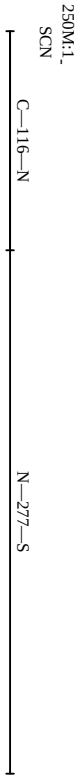
Molecular Origami of SCN^-

!C1
N1
S1
SCN^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.06 offsety 0.9

View -1



Current: (centerx 4.24) (centery 5.90) (scale 250)
%%BoundingBox: 295 136 315 435 actual: 305 146 305 425 center: 305 286 actual size: 0 278
Better: (centerx 4.25) (centery 7.43) (scale 250)
%%BoundingBox: 292 312 312 610 actual: 302 322 302 600 center: 302 461 actual size: 0 278