Molecular Origami of SCN^-

given information

ElementNames	[(C) (N) (S)]	
dotted	F	
distance	115.628	C^1-N^1
distance	164.924	C^1-S^1
angle	179.034	S^1-C^1-N^1
	280.5	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 -0.07 offsety 0.9

View -1

Current: (centerx 4.23) (centery 5.90) (scale 250)
%%BoundingBox: 295 134 315 435 actual: 305 144 305 425 center: 305 284 actual size: 0 281
Better: (centerx 4.25) (centery 7.45) (scale 250)
%%BoundingBox: 291 310 311 611 actual: 301 320 301 601 center: 301 461 actual size: 0 281