

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

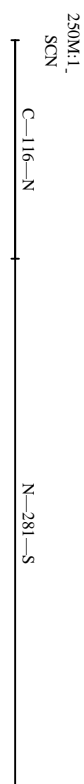
Molecular Origami of SCN⁻

!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