Molecular Origami of C2H2

given information

ElementNames	[(C) (C) (H)]	
dotted	F	
distance	101.533	C^1-H^1
distance	118.645	C^1-C^2
angle	180.000	H^1-C^1-C^2
	220.2	H^1-C^2
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

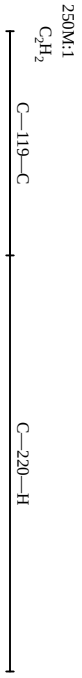
Molecular Origami of C2H2

!C1
C2
H1
C2H2

linear

scale 250,000,000 : 1
units: pm
offsetx -0.05 offsety 0.9

View -1



Current: (centerx 4.25) (centery 5.90) (scale 250)
%%BoundingBox: 296 175 316 435 actual: 306 185 306 425 center: 306 305 actual size: 0 240
Better: (centerx 4.25) (centery 7.17) (scale 250)
%%BoundingBox: 292 331 312 591 actual: 302 341 302 581 center: 302 461 actual size: 0 240