Molecular Origami of Se(CH3)2

given information

ElementNames	[(Se) (C) (C)]	
dotted	F	
distance	198.000	Se^1-C^1
distance	198.033	Se^1-C^2
angle	98.011	C^2-Se^1-C^1
	298.9	C^2-C^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

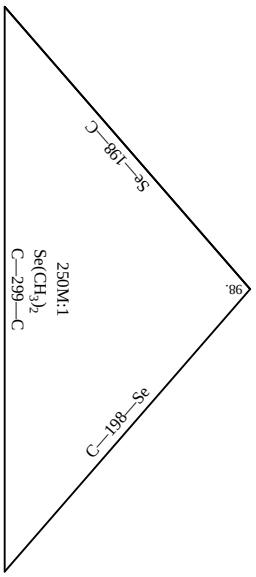
Molecular Origami of Se(CH3)2

!Se1
C1
C2
Se(CH3)2

bent

scale 250,000,000 : 1
units: pm
offsetx -1.51 offsety 0.9

View -1



Current: (centerx 2.79) (centery 5.90) (scale 250)
%%BoundingBox: 191 309 303 541 actual: 201 319 293 531 center: 247 425 actual size: 92 212
Better: (centerx 3.61) (centery 5.50) (scale 250)
%%BoundingBox: 141 345 253 577 actual: 151 355 243 567 center: 197 461 actual size: 92 212