Molecular Origami of O(CH3)2

given information

ElementNames	[(O) (C) (C)]	
dotted	F	
distance	135.439	O^1-C^1
distance	148.133	O^1-C^2
angle	110.445	C^2-O^1-C^1
	233.	C^2-C^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

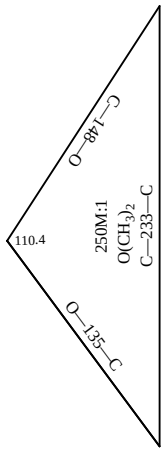
Molecular Origami of O(CH3)2

!O1
C1
C2
O(CH3)2

bent

scale 250,000,000 : 1
units: pm
offsetx 1.0 offsety 0.9

View -1



Current: (centerx 5.30) (centery 5.90) (scale 250)
%%BoundingBox: 314 332 392 517 actual: 324 342 382 507 center: 353 425 actual size: 57 165
Better: (centerx 4.65) (centery 5.50) (scale 250)
%%BoundingBox: 339 368 417 553 actual: 349 378 407 543 center: 378 461 actual size: 57 165