

Molecular Origami of S(CH₃)₂

given information

ElementNames	[(S) (C) (C)]	
dotted	F	
distance	180.000	S ¹ -C ¹
distance	180.007	S ¹ -C ²
angle	98.981	C ² -S ¹ -C ¹
	273.7	C ² -C ¹
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

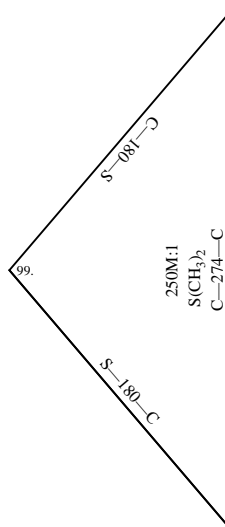
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!S1
C1
C2
S(CH₃)₂

bent

scale 250,000,000 : 1
units: pm
offsetx 1.28 offsety 0.9

View -1



Current: (centerx 5.58) (centery 5.90) (scale 250)

%%BoundingBox: 309 318 412 532 actual: 319 328 402 522

center: 360 425

actual size: 83 194

Better: (centerx 4.83) (centery 5.50) (scale 250)

%%BoundingBox: 347 354 450 568

actual: 357 364 440 558

center: 398 461

actual size: 83 194