

Molecular Origami of CO<sub>3</sub><sup>2-</sup>-  
given information

ElementNames	[ (C) (O) (O) (O) ]	
dotted	F	
distance	127.161	C <sup>1</sup> -O <sup>2</sup>
distance	127.186	C <sup>1</sup> -O <sup>1</sup>
distance	127.256	C <sup>1</sup> -O <sup>3</sup>
angle	119.967	O <sup>3</sup> -C <sup>1</sup> -O <sup>1</sup>
	220.3	O <sup>3</sup> -O <sup>1</sup>
angle	119.984	O <sup>3</sup> -C <sup>1</sup> -O <sup>2</sup>
	220.3	O <sup>3</sup> -O <sup>2</sup>
angle	120.036	O <sup>2</sup> -C <sup>1</sup> -O <sup>1</sup>
	220.3	O <sup>2</sup> -O <sup>1</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

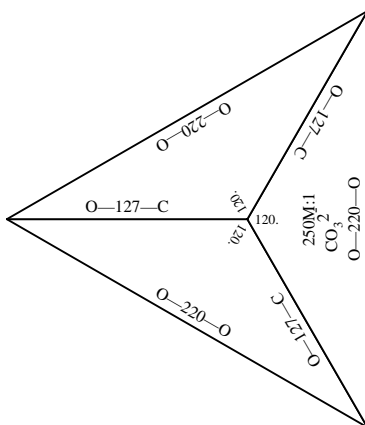
Molecular Origami of CO<sub>3</sub><sup>2-</sup>

!C1  
O1  
O2  
O3  
CO<sub>3</sub><sup>2-</sup>

trigonal planar

scale 250,000,000 : 1  
units: pm  
offsetx 0.89 offsety 0.9

View -1



Current: (centerx 5.19) (centery 5.90) (scale 250)

%%BoundingBox: 228 337 384 513 actual: 238 347 374 503

center: 306 425

actual size: 135 156

Better: (centerx 5.19) (centery 5.50) (scale 250)

%%BoundingBox: 292 373 448 549

actual: 302 383 438 539

center: 370 461

actual size: 135 156