

Molecular Origami of C2O4<sup>2-</sup>-  
given information

ElementNames	[ (C) (C) (O) (O) ]	
dotted	F	
distance	120.322	C <sup>1</sup> -O <sup>4</sup>
distance	132.855	C <sup>1</sup> -O <sup>1</sup>
distance	152.115	C <sup>1</sup> -C <sup>2</sup>
angle	114.855	O <sup>1</sup> -C <sup>1</sup> -C <sup>2</sup>
	240.4	O <sup>1</sup> -C <sup>2</sup>
angle	121.916	O <sup>4</sup> -C <sup>1</sup> -O <sup>1</sup>
	221.4	O <sup>4</sup> -O <sup>1</sup>
angle	123.221	O <sup>4</sup> -C <sup>1</sup> -C <sup>2</sup>
	240.1	O <sup>4</sup> -C <sup>2</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

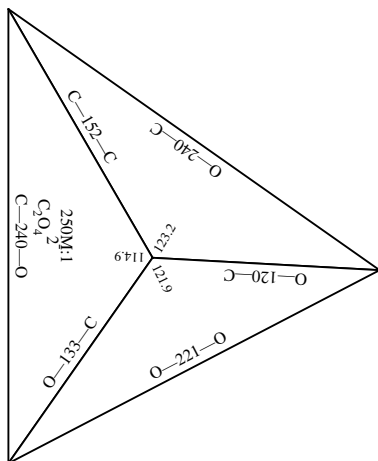
Molecular Origami of C2O4^2-

!C1  
C2  
O1  
O4  
C2O4^2-

trigonal planar

scale 250,000,000 : 1  
units: pm  
offsetx -1.1 offsety 0.9

View -1



Current: (centerx 3.20) (centery 5.90) (scale 250)

%%BoundingBox: 220 330 380 520 actual: 230 340 370 510

Better: (centerx 3.28) (centery 5.50) (scale 250)

%%BoundingBox: 147 366 306 556 actual: 157 376 296 546

center: 300 425

actual size: 139 170

center: 227 461

actual size: 139 170