

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

ElementNames	[ (C) (C) (O) (O) ]	
dotted	F	
distance	123.592	C <sup>1</sup> -O <sup>1</sup>
distance	125.771	C <sup>1</sup> -O <sup>2</sup>
distance	154.795	C <sup>1</sup> -C <sup>2</sup>
angle	114.357	O <sup>2</sup> -C <sup>1</sup> -C <sup>2</sup>
	236.3	O <sup>2</sup> -C <sup>2</sup>
angle	118.467	O <sup>1</sup> -C <sup>1</sup> -C <sup>2</sup>
	239.7	O <sup>1</sup> -C <sup>2</sup>
angle	127.177	O <sup>2</sup> -C <sup>1</sup> -O <sup>1</sup>
	223.3	O <sup>2</sup> -O <sup>1</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

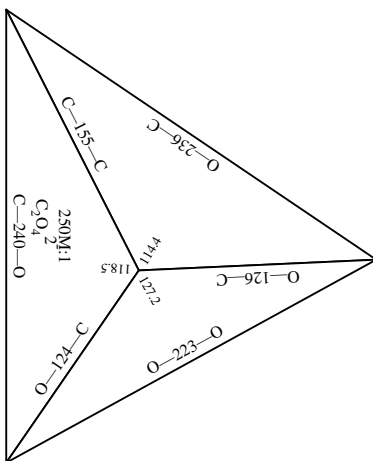
Molecular Origami of C2O4^2-

!C1  
C2  
O1  
O2  
C2O4^2-

trigonal planar

scale 250,000,000 : 1  
units: pm  
offsetx -1.08 offsety 0.9

View -1



Current: (centerx 3.22) (centery 5.90) (scale 250)

%%BoundingBox: 222 330 381 520 actual: 232 340 371 510

center: 301 425

actual size: 139 170

Better: (centerx 3.29) (centery 5.50) (scale 250)

%%BoundingBox: 149 366 308 556

actual: 159 376 298 546

center: 228 461

actual size: 139 170