

Molecular Origami of  $\text{HCO}_3^-$ -  
given information

ElementNames	[ (C) (O) (O) (O) ]	
dotted	F	
distance	123.212	$\text{C}^1\text{-O}^2$
distance	125.733	$\text{C}^1\text{-O}^1$
distance	133.019	$\text{C}^1\text{-O}^3$
angle	117.052	$\text{O}^3\text{-C}^1\text{-O}^2$
	218.6	$\text{O}^3\text{-O}^2$
angle	117.199	$\text{O}^3\text{-C}^1\text{-O}^1$
	220.9	$\text{O}^3\text{-O}^1$
angle	125.734	$\text{O}^2\text{-C}^1\text{-O}^1$
	221.6	$\text{O}^2\text{-O}^1$
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

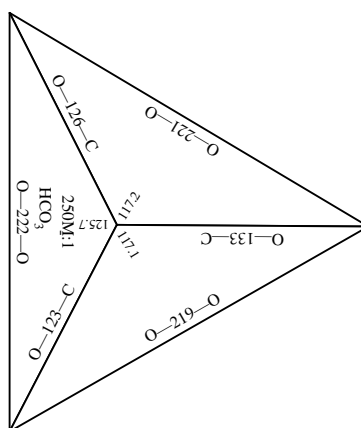
Molecular Origami of  $\text{HCO}_3^-$

!C1  
O1  
O2  
O3  
 $\text{HCO}_3^-$

trigonal planar

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

View -1



Current: (centerx 3.30) (centery 5.90) (scale 250)

%%BoundingBox: 228 336 382 513 actual: 238 346 372 503

center: 305 425

actual size: 134 157

Better: (centerx 3.32) (centery 5.50) (scale 250)

%%BoundingBox: 157 372 311 549

actual: 167 382 301 539

center: 234 461

actual size: 134 157