

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

ElementNames	[ (C) (O) (O) (O) ]	
dotted	F	
distance	122.757	$\text{C}^1\text{-O}^2$
distance	126.443	$\text{C}^1\text{-O}^1$
distance	131.317	$\text{C}^1\text{-O}^3$
angle	117.918	$\text{O}^3\text{-C}^1\text{-O}^2$
	217.7	$\text{O}^3\text{-O}^2$
angle	118.856	$\text{O}^3\text{-C}^1\text{-O}^1$
	221.9	$\text{O}^3\text{-O}^1$
angle	123.226	$\text{O}^2\text{-C}^1\text{-O}^1$
	219.2	$\text{O}^2\text{-O}^1$
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

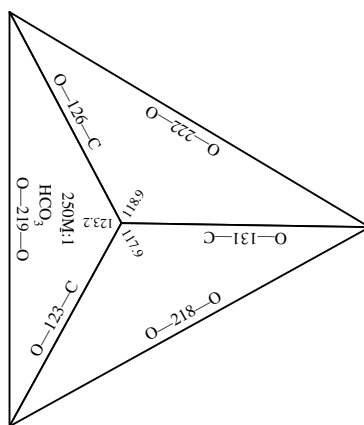
Molecular Origami of  $\text{HCO}_3^-$

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

trigonal planar

scale 250,000,000 : 1  
units: pm  
offsetx -0.99 offsety 0.9

View -1



Current: (centerx 3.31) (centery 5.90) (scale 250)

%%BoundingBox: 228 337 383 512 actual: 238 347 373 502

center: 306 425

actual size: 135 155

Better: (centerx 3.31) (centery 5.50) (scale 250)

%%BoundingBox: 157 373 312 548

actual: 167 383 302 538

center: 235 461

actual size: 135 155