

Molecular Origami of HCO_3^- -
given information

ElementNames	[(C) (O) (O) (O)]	
dotted	F	
distance	123.004	$\text{C}^1\text{-O}^3$
distance	125.340	$\text{C}^1\text{-O}^1$
distance	126.206	$\text{C}^1\text{-O}^2$
angle	115.542	$\text{O}^2\text{-C}^1\text{-O}^1$
	212.8	$\text{O}^2\text{-O}^1$
angle	119.862	$\text{O}^3\text{-C}^1\text{-O}^1$
	214.9	$\text{O}^3\text{-O}^1$
angle	124.592	$\text{O}^3\text{-C}^1\text{-O}^2$
	220.6	$\text{O}^3\text{-O}^2$
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

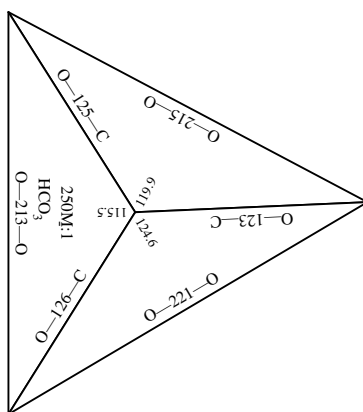
Molecular Origami of HCO_3^-

!C1
O1
O2
O3
 HCO_3^-

trigonal planar

scale 250,000,000 : 1
units: pm
offsetx -0.98 offsety 0.9

View -1



Current: (centerx 3.32) (centery 5.90) (scale 250)

%%BoundingBox: 229 339 384 510 actual: 239 349 374 500

center: 306 425

actual size: 135 151

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

%%BoundingBox: 158 375 313 546

actual: 168 385 303 536

center: 235 461

actual size: 135 151