

# Molecular Origami of CH<sub>3</sub>CO<sub>2</sub><sup>-</sup>

given information

ElementNames	[ (C) (C) (O) (O) ]	
dotted	F	
distance	124.348	C <sup>1</sup> -O <sup>1</sup>
distance	124.425	C <sup>1</sup> -O <sup>2</sup>
distance	151.865	C <sup>1</sup> -C <sup>2</sup>
angle	117.644	O <sup>1</sup> -C <sup>1</sup> -C <sup>2</sup>
	236.7	O <sup>1</sup> -C <sup>2</sup>
angle	118.440	O <sup>2</sup> -C <sup>1</sup> -C <sup>2</sup>
	237.8	O <sup>2</sup> -C <sup>2</sup>
angle	123.916	O <sup>2</sup> -C <sup>1</sup> -O <sup>1</sup>
	219.6	O <sup>2</sup> -O <sup>1</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

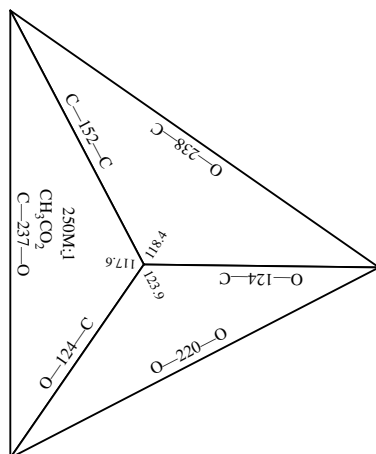
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!C1  
C2  
O1  
O2  
CH<sub>3</sub>CO<sub>2</sub><sup>-</sup>

trigonal planar

scale 250,000,000 : 1  
units: pm  
offsetx -1.09 offsety 0.9

View -1



Current: (centerx 3.21) (centery 5.90) (scale 250)

%%BoundingBox: 221 331 379 519 actual: 231 341 369 509

center: 300 425

actual size: 138 168

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

%%BoundingBox: 148 367 307 555

actual: 158 377 297 545

center: 228 461

actual size: 138 168