Molecular Origami of CH2Br2

given information

ElementNames	[(C) (Br) (Br) (H) (H)]	
distance	99.999	C^1-H^2
distance	100.007	C^1-H^1
distance	188.199	C^1-Br^2
distance	190.256	C^1-Br^1
angle	108.479	H^2-C^1-Br^2
	239.5	H^2-Br^2
angle	108.481	H^1-C^1-Br^2
	239.5	H^1-Br^2
angle	108.482	H^1-C^1-Br^1
	241.4	H^1-Br^1
angle	108.482	H^2-C^1-Br^1
	241.4	H^2-Br^1
angle	109.501	H^2-C^1-H^1
	163.3	H^2-H^1
angle	113.370	Br^2-C^1-Br^1
	316.3	Br^2-Br^1
dopage	T	
AutoAlign	F	

structure type: XABCD

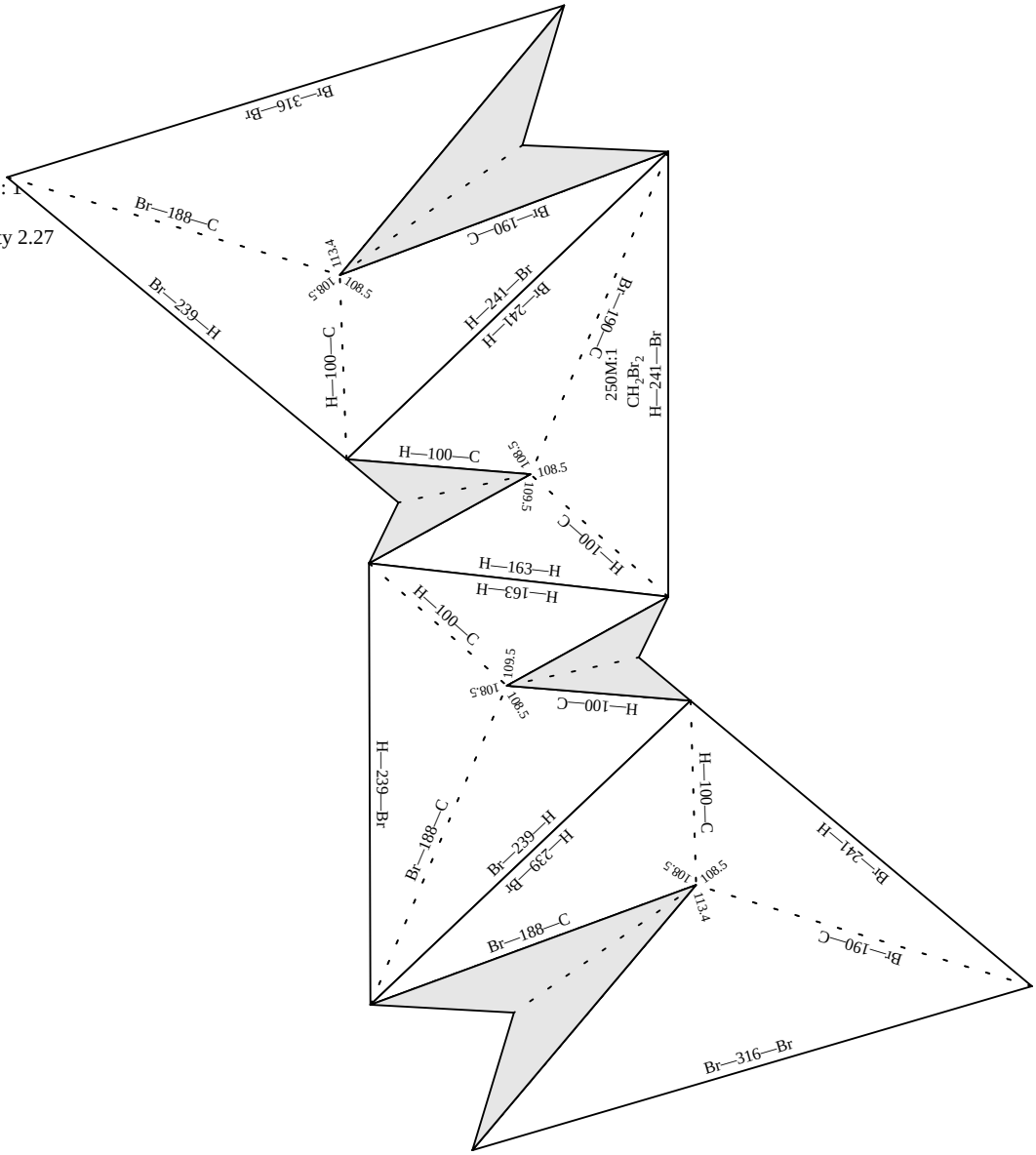
Molecular Origami of CH2Br2

!C1
Br1
Br2
H1
H2
CH2Br2

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx 0.36 offsety 2.27

View -1



Current: (centerx 4.66) (centery 7.27) (scale 250)
%%BoundingBox: 124 254 538 714 actual: 134 264 528 704 center: 331 484 actual size: 394 440
Better: (centerx 4.31) (centery 6.05) (scale 250)
%%BoundingBox: 125 329 539 789 actual: 135 339 529 779 center: 332 559 actual size: 394 440