Molecular Origami of CHBr3

given information

ElementNames	[(C) (Br) (Br) (Br) (H)]	
distance	100.002	C^1-H^1
distance	191.225	C^1-Br^3
distance	191.225	C^1-Br^1
distance	195.015	C^1-Br^2
angle	106.522	H^1-C^1-Br^3
	239.7	H^1-Br^3
angle	106.522	H^1-C^1-Br^1
	239.7	H^1-Br^1
angle	106.892	H^1-C^1-Br^2
	243.6	H^1-Br^2
angle	112.033	Br^3-C^1-Br^2
	320.3	Br^3-Br^2
angle	112.033	Br^2-C^1-Br^1
	320.3	Br^2-Br^1
angle	112.362	Br^3-C^1-Br^1
	317.7	Br^3-Br^1
dopage	T	
AutoAlign	F	

structure type: XABCD

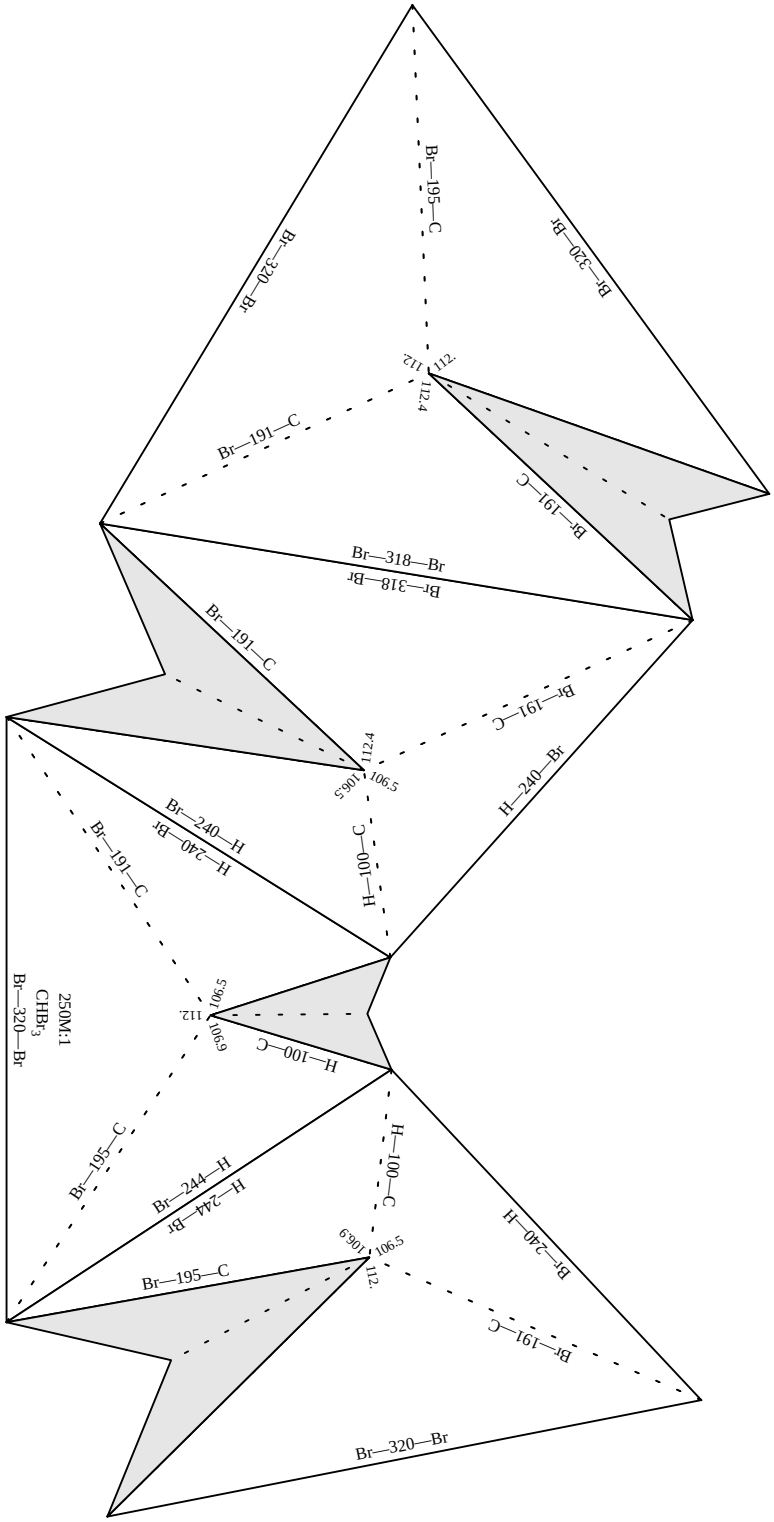
Molecular Origami of CHBr3

!C1
Br1
Br2
Br3
H1
CHBr3

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.99 offsety -0.46

View -1



Current: (centerx 3.31) (centery 4.54) (scale 250)
%%BoundingBox: 152 129 458 716 actual: 162 139 448 706 center: 305 422 actual size: 286 567
Better: (centerx 3.33) (centery 4.17) (scale 250)
%%BoundingBox: 82 69 388 656 actual: 92 79 378 646 center: 235 363 actual size: 286 567