Molecular Origami of CHBr3

given information

ElementNames	[(C) (Br) (Br) (Br) (H)]	
distance	96.934	C^1-H^1
distance	188.965	C^1-Br^2
distance	193.828	C^1-Br^3
distance	196.332	C^1-Br^1
angle	108.074	Br^3-C^1-Br^1
	315.8	Br^3-Br^1
angle	108.770	H^1-C^1-Br^2
	238.5	H^1-Br^2
angle	109.048	Br^2-C^1-Br^1
	313.8	Br^2-Br^1
angle	109.629	H^1-C^1-Br^3
	244.1	H^1-Br^3
angle	110.389	Br^3-C^1-Br^2
	314.3	Br^3-Br^2
angle	110.928	H^1-C^1-Br^1
	248.1	H^1-Br^1
dopage	T	
AutoAlign	F	

structure type: XABCD

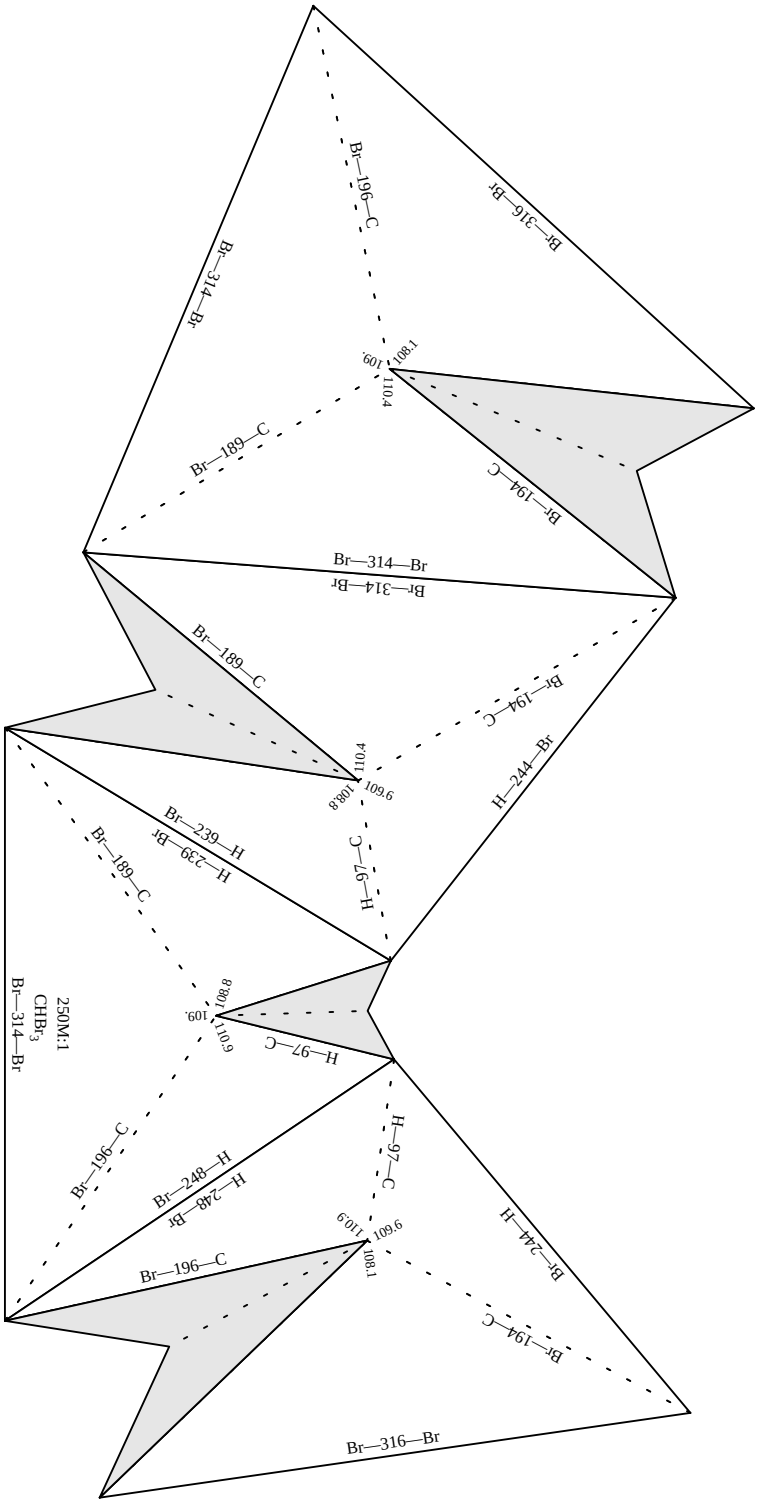
Molecular Origami of CHBr3

!C1
Br1
Br2
Br3
H1
CHBr3

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.94 offsety -0.53

View -1



Current: (centerx 3.36) (centery 4.47) (scale 250)
%%BoundingBox: 153 131 454 711 actual: 163 141 444 701 center: 303 421 actual size: 281 559
Better: (centerx 3.40) (centery 4.12) (scale 250)
%%BoundingBox: 88 68 389 648 actual: 98 78 379 638 center: 238 358 actual size: 281 559