Molecular Origami of CBr4

given information

ElementNames	[(C) (Br) (Br) (Br) (Br)]	
distance	182.492	C^1-Br^4
distance	190.167	C^1-Br^2
distance	193.385	C^1-Br^1
distance	195.430	C^1-Br^3
angle	105.429	Br^4-C^1-Br^2
	296.5	Br^4-Br^2
angle	107.424	Br^3-C^1-Br^1
	313.4	Br^3-Br^1
angle	109.511	Br^4-C^1-Br^3
	308.7	Br^4-Br^3
angle	110.086	Br^2-C^1-Br^1
	314.4	Br^2-Br^1
angle	111.102	Br^4-C^1-Br^1
	310.	Br^4-Br^1
angle	113.335	Br^3-C^1-Br^2
	322.2	Br^3-Br^2
dopage	T	
AutoAlign	F	

structure type: XABCD

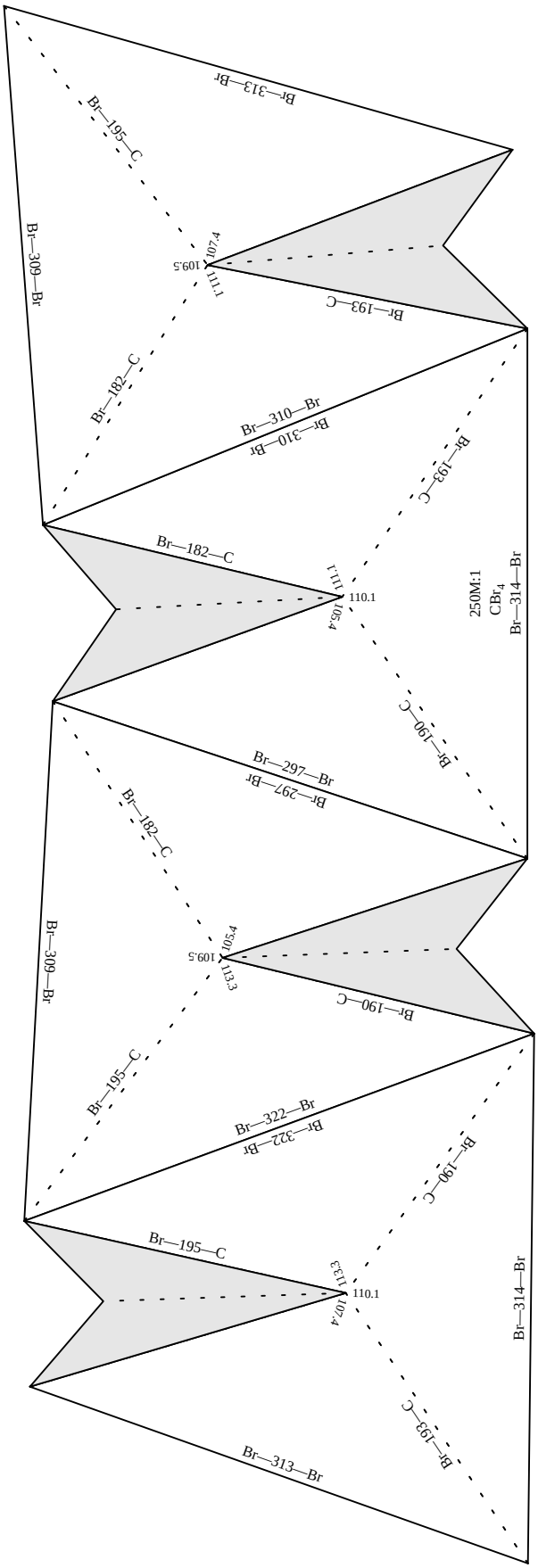
Molecular Origami of CBr4

!C1
Br1
Br2
Br3
Br4
CBr4

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx 0.39 offsety 2.02

View -1



Current: (centerx 4.69) (centery 7.02) (scale 250)
%%BoundingBox: 186 89 428 764 actual: 196 99 418 754 center: 307 426 actual size: 223 655
Better: (centerx 4.68) (centery 6.60) (scale 250)
%%BoundingBox: 213 204 455 879 actual: 223 214 445 869 center: 334 541 actual size: 223 655