Molecular Origami of CH3I

given information

ElementNames	[(C) (H) (H) (H) (I)]	
distance	99.993	C^1-H^1
distance	100.002	C^1-H^2
distance	100.007	C^1-H^3
distance	214.865	C^1-I^1
angle	109.440	H^3-C^1-H^1
	163.3	H^3-H^1
angle	109.442	H^2-C^1-H^1
	163.3	H^2-H^1
angle	109.444	H^3-C^1-H^2
	163.3	H^3-H^2
angle	109.499	I^1-C^1-H^3
	265.5	I^1-H^3
angle	109.499	I^1-C^1-H^2
	265.5	I^1-H^2
angle	109.504	I^1-C^1-H^1
	265.5	I^1-H^1
dopage	T	
AutoAlign	F	

structure type: XABCD

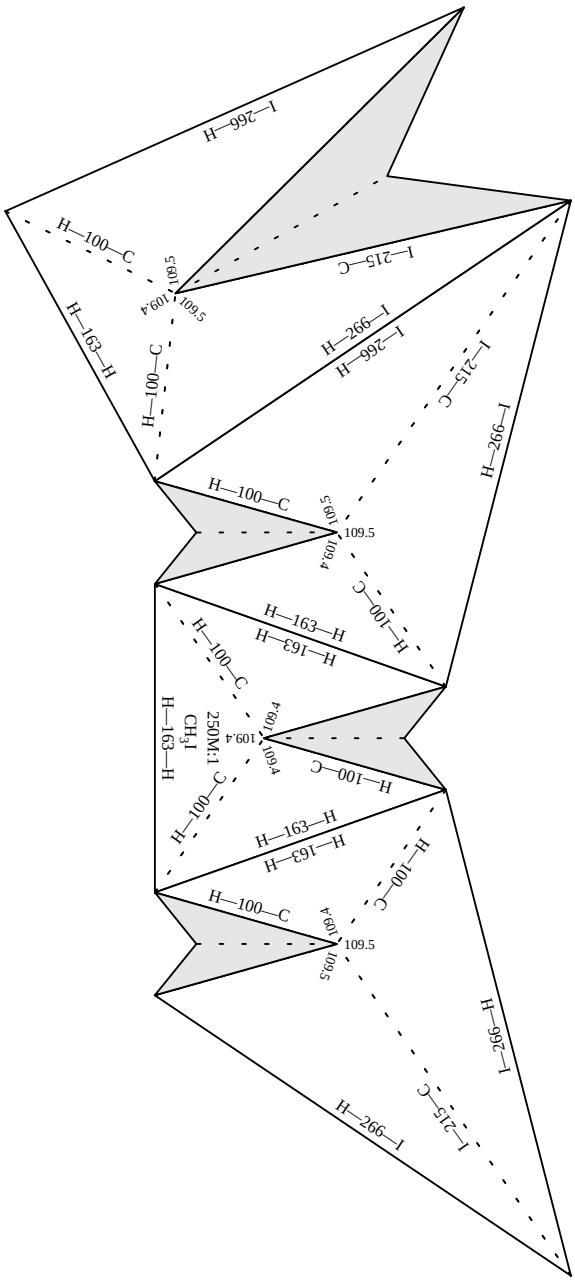
Molecular Origami of CH3I

!C1
H1
H2
H3
I1
CH3I

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.18 offsety 0.4

View -1



Current: (centerx 4.12) (centery 5.40) (scale 250)
%%BoundingBox: 190 177 422 673 actual: 200 187 412 663 center: 306 425 actual size: 212 476
Better: (centerx 4.12) (centery 5.00) (scale 250)
%%BoundingBox: 177 177 409 673 actual: 187 187 399 663 center: 293 425 actual size: 212 476