Molecular Origami of CH2F2

given information

ElementNames	[(C) (F) (F) (H) (H)]	
distance	105.966	C^1-H^1
distance	107.221	C^1-H^2
distance	134.353	C^1-F^2
distance	137.110	C^1-F^1
angle	103.557	H^1-C^1-F^1
	191.9	H^1-F^1
angle	104.471	H^2-C^1-F^1
	194.	H^2-F^1
angle	104.499	H^1-C^1-F^2
	190.8	H^1-F^2
angle	105.125	F^2-C^1-F^1
	215.6	F^2-F^1
angle	110.378	H^2-C^1-F^2
	198.9	H^2-F^2
angle	126.839	H^2-C^1-H^1
	190.7	H^2-H^1
dopage	T	
AutoAlign	F	

structure type: XABCD

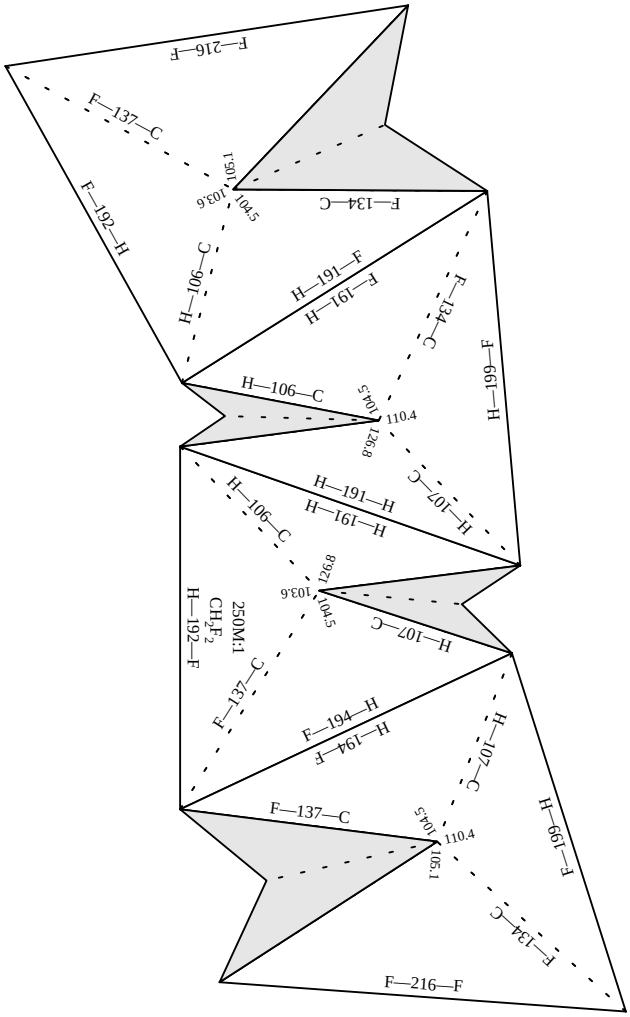
Molecular Origami of CH2F2

!C1
F1
F2
H1
H2
CH2F2

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.2 offsety 0.21

View -1



Current: (centerx 4.10) (centery 5.21) (scale 250)
%%BoundingBox: 168 207 420 605 actual: 178 217 410 595 center: 294 406 actual size: 233 377
Better: (centerx 4.27) (centery 5.07) (scale 250)
%%BoundingBox: 165 212 418 610 actual: 175 222 408 600 center: 292 411 actual size: 233 377