

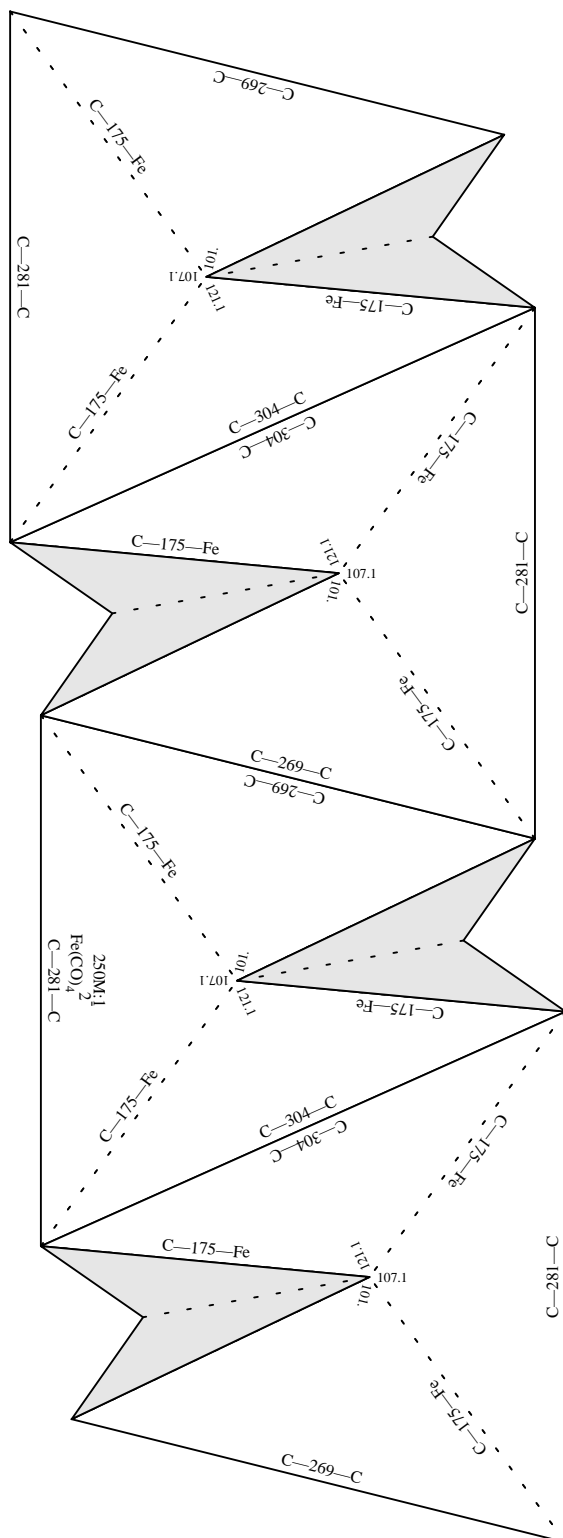
Molecular Origami of Fe(CO)₄²⁻
given information

ElementNames	[(Fe) (C) (C) (C) (C)]	
distance	174.649	Fe ¹ -C ³
distance	174.649	Fe ¹ -C ²
distance	174.655	Fe ¹ -C ⁴
distance	174.655	Fe ¹ -C ¹
angle	100.961	C ⁴ -Fe ¹ -C ¹
	269.5	C ⁴ -C ¹
angle	100.966	C ³ -Fe ¹ -C ²
	269.5	C ³ -C ²
angle	107.055	C ⁴ -Fe ¹ -C ³
	280.9	C ⁴ -C ³
angle	107.055	C ² -Fe ¹ -C ¹
	280.9	C ² -C ¹
angle	121.100	C ⁴ -Fe ¹ -C ²
	304.2	C ⁴ -C ²
angle	121.100	C ³ -Fe ¹ -C ¹
	304.2	C ³ -C ¹
dopage	T	
AutoAlign	F	

structure type: XABCD

!Fe1
C1
C2
C3
C4
Fe(CO)4²⁻

```
scale 250,000,000 : 1
units: pm
offsetx -0.31 offsety -0.16
View -1
```



actual size: 208 574

actual size: 208 574