

Molecular Origami of Mo(pyrazoylborate)Cl₃facial
given information

ElementNames	[(Mo) (Cl) (Cl) (Cl) (N) (N) (N)]	
distance	218.116	Mo ¹ -N ⁵
distance	218.129	Mo ¹ -N ²
distance	219.760	Mo ¹ -N ¹
distance	243.042	Mo ¹ -Cl ³
distance	243.049	Mo ¹ -Cl ²
distance	243.307	Mo ¹ -Cl ¹
angle	85.686	N ² -Mo ¹ -N ¹
	297.8	N ² -N ¹
angle	85.688	N ⁵ -Mo ¹ -N ¹
	297.8	N ⁵ -N ¹
angle	86.797	N ⁵ -Mo ¹ -N ²
	299.7	N ⁵ -N ²
angle	89.769	N ⁵ -Mo ¹ -Cl ¹
	326.1	N ⁵ -Cl ¹
angle	89.771	N ² -Mo ¹ -Cl ¹
	326.1	N ² -Cl ¹
angle	90.873	N ² -Mo ¹ -Cl ²
	329.	N ² -Cl ²
angle	90.875	N ⁵ -Mo ¹ -Cl ³
	329.	N ⁵ -Cl ³
angle	91.275	N ¹ -Mo ¹ -Cl ²
	331.3	N ¹ -Cl ²
angle	91.275	N ¹ -Mo ¹ -Cl ³
	331.3	N ¹ -Cl ³
angle	91.305	Cl ³ -Mo ¹ -Cl ²
	347.6	Cl ³ -Cl ²
angle	93.097	Cl ² -Mo ¹ -Cl ¹
	353.1	Cl ² -Cl ¹
angle	93.097	Cl ³ -Mo ¹ -Cl ¹
	353.1	Cl ³ -Cl ¹
angle	173.743	N ¹ -Mo ¹ -Cl ¹
	462.4	N ¹ -Cl ¹
angle	176.300	N ⁵ -Mo ¹ -Cl ²
	460.9	N ⁵ -Cl ²
angle	176.300	N ² -Mo ¹ -Cl ³
	460.9	N ² -Cl ³
dopage	T	
AutoAlign	F	

structure type: XABCDEF

!Mo1
Cl1
Cl2
Cl3
N1
N2
N5
Mo(pyrazoylborate)Cl3facial

```
scale 125,000,000 : 1
units: pm
offsetx -0.01 offsety 0.08
```

[illegible]

actual size: 343 710

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