

(23) **Today**

Section 6.4.6: Halogen Bonds

Section 9.1 - 9.3

Introduction to Coordination Chemistry,
Nomenclature and Ligands

Next Class (24)

Section 9.3 – 9.5

Nomenclature and Ligands, Isomerism,
Coordination Number and Structures

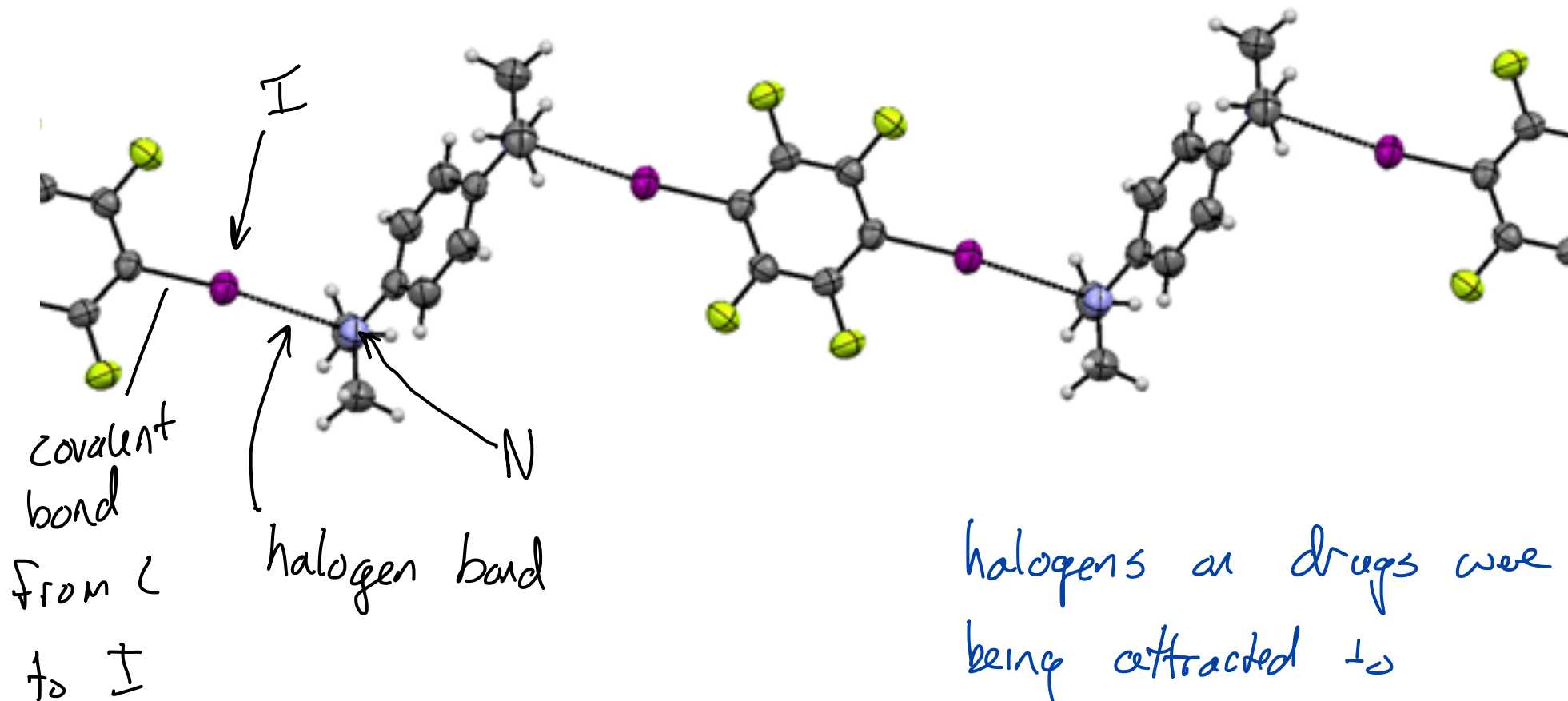
(25) **Second Class from Today**

9.5 Coordination Number and Structures

Chap 10

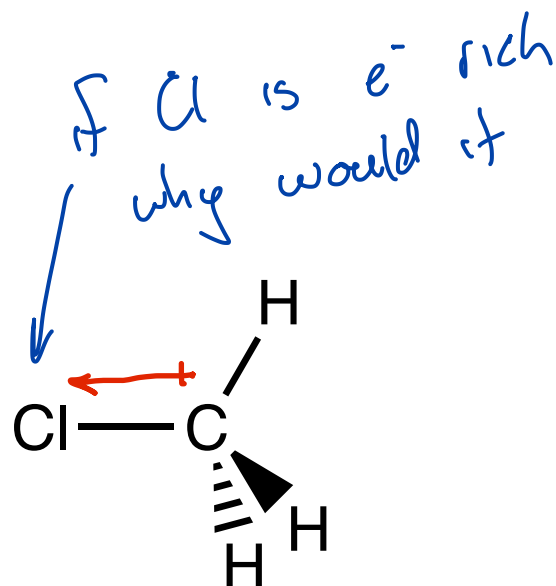
Third Class from Today (26)

Test 2

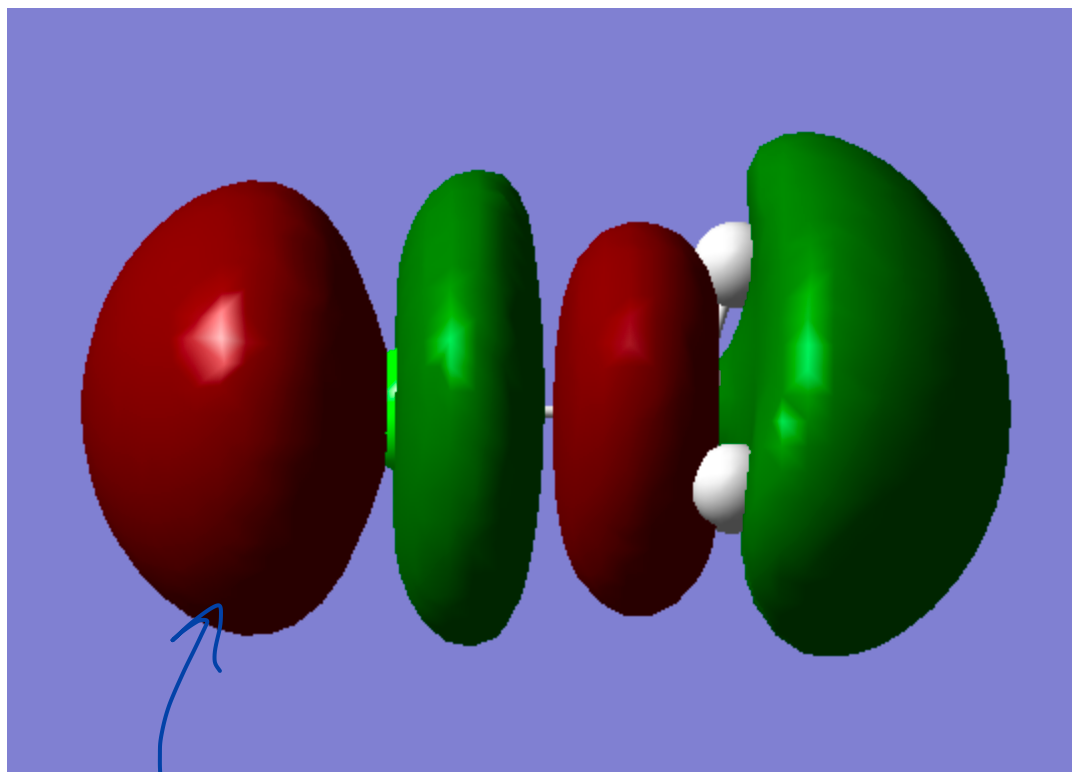


halogens in drugs were
being attracted to
C=O O atoms in peptides

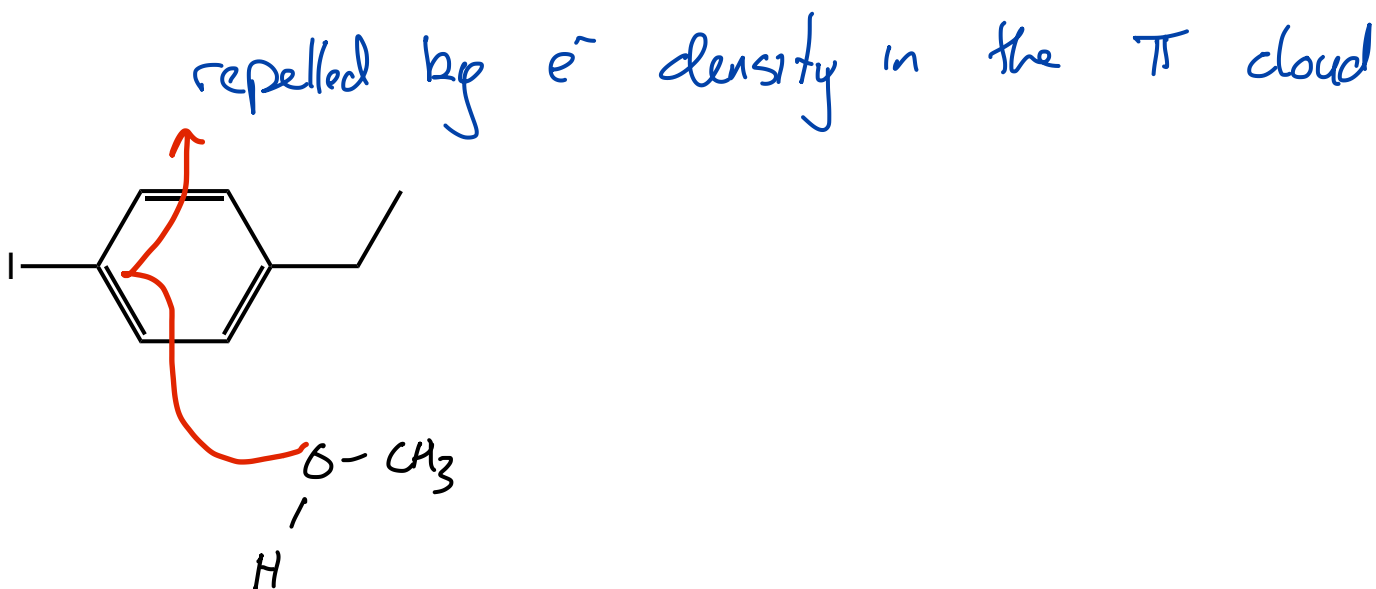
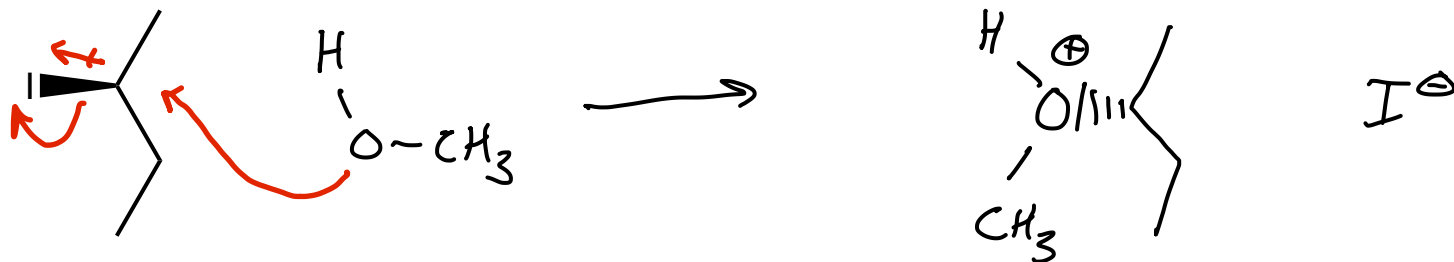
[https://chem.libretexts.org/Bookshelves/Inorganic_Chemistry/Inorganic_Chemistry_\(LibreTexts\)/06: Acid-Base and Donor-Acceptor Chemistry/6.04: Lewis Concept and Frontier Orbitals/6.4.06: Lewis base strength may also be estimated by measuring structural or energy changes on formation of a Lewis acid-base complex as illustrated by efforts to spectroscopically assess of the strengths of halogen bonds](https://chem.libretexts.org/Bookshelves/Inorganic_Chemistry/Inorganic_Chemistry_(LibreTexts)/06:_Acid-Base_and_Donor-Acceptor_Chemistry/6.04:_Lewis_Concept_and_Frontier_Orbitals/6.4.06:_Lewis_base_strength_may_also_be_estimated_by_measuring_structural_or_energy_changes_on_formation_of_a_Lewis_acid-base_complex_as_illustrated_by_efforts_to_spectroscopically_assess_of_the_strengths_of_halogen_bonds)
The figure was produced using Mercury 3.5.1 based on Cambridge crystallographic database entry 172272. This work by Stephen Contakes is licensed under a Creative Commons Attribution 4.0 International License.

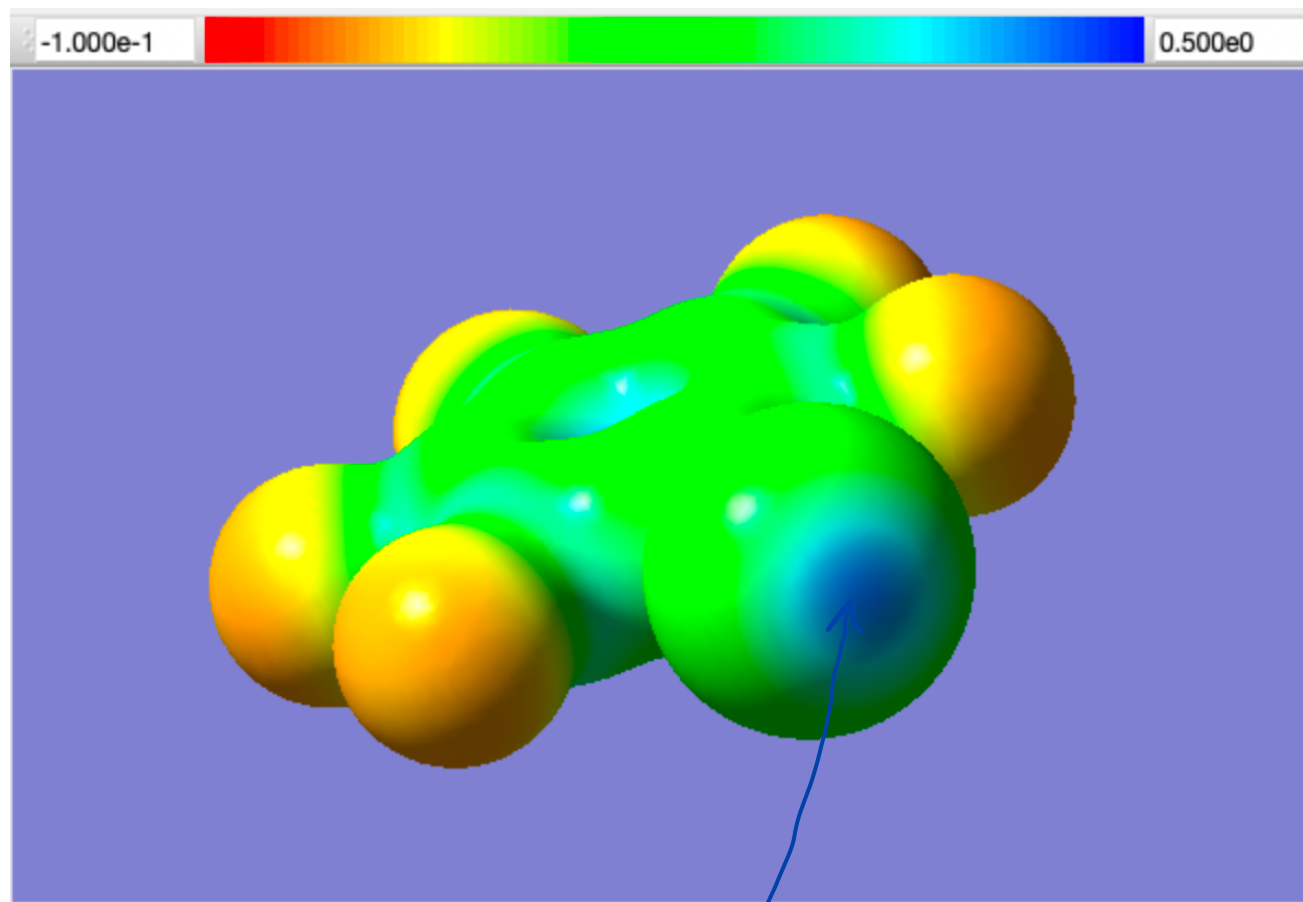
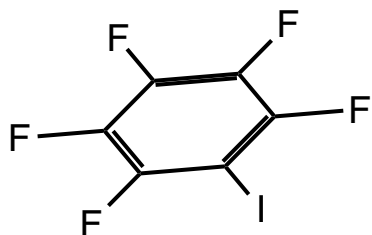


LUMO



the Cl end of the LUMO can
accept e^- density, too.

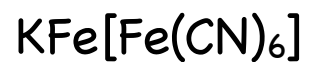




δ^+ spot on the I

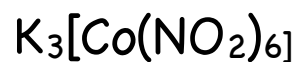
a σ -hole so e^- rich atoms
can be attracted to the σ -hole⁴⁵

Prussian Blue



https://en.wikipedia.org/wiki/Prussian_blue

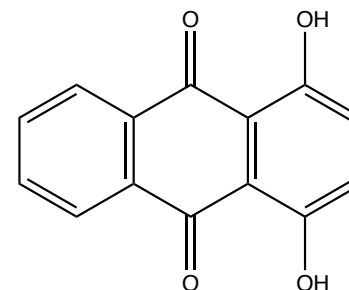
Aureolin



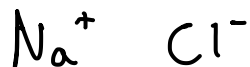
https://en.wikipedia.org/wiki/Potassium_cobaltinitrite

Alizarin

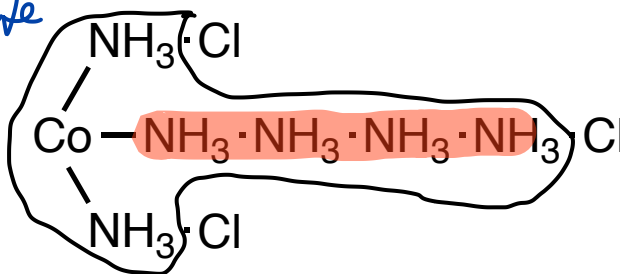
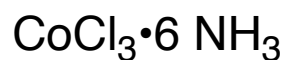
Ca^{2+} salts of



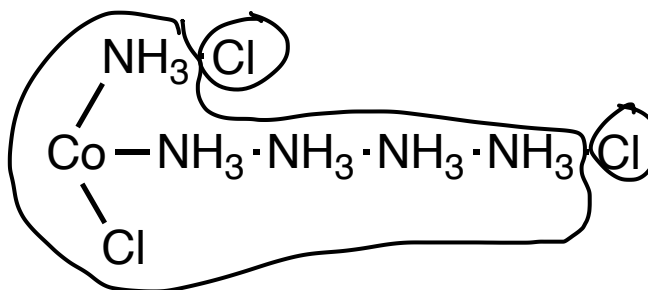
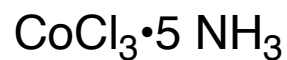
<https://en.wikipedia.org/wiki/Alizarin>



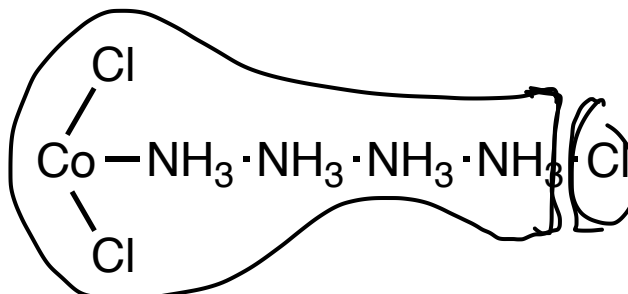
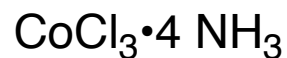
valence = oxidation state



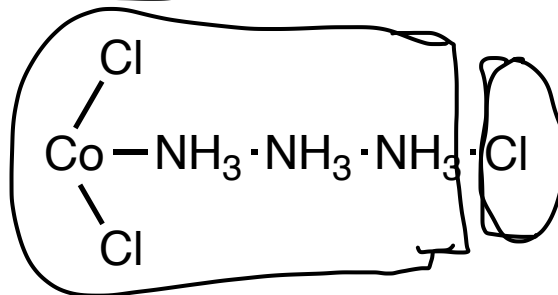
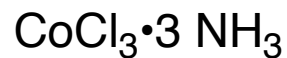
4 ions



3 ions



2 ions



2 ions

Cl⁻ ions
precipitated by
Ag⁺

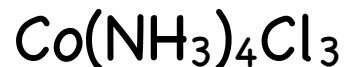
PtCl₄•6NH₃ 4

PtCl₄•5NH₃ 3

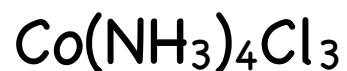
PtCl₄•4NH₃ 2

PtCl₄•3NH₃ 1

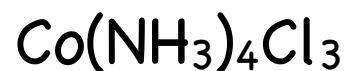
PtCl₄•2NH₃ 0



two forms of $\text{Co}(\text{NH}_3)_4\text{Cl}_3$



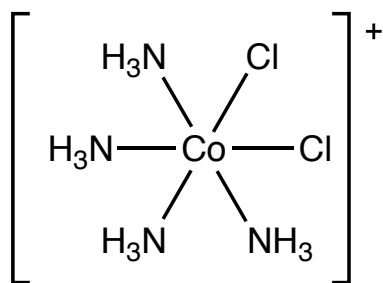
purple



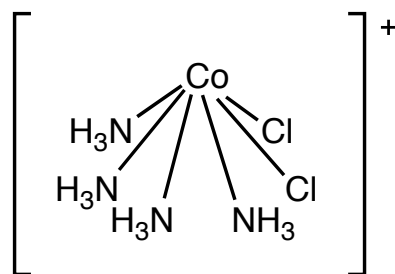
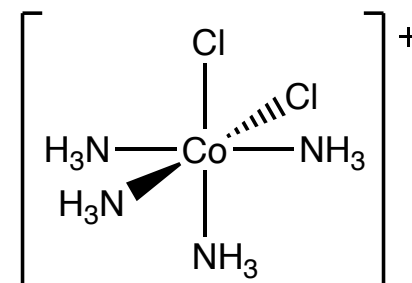
green

same chemical formulas but
different structures 'cuz
one version is green & the
other is purple

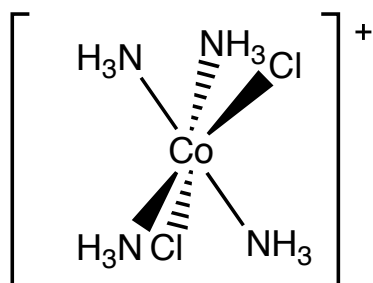
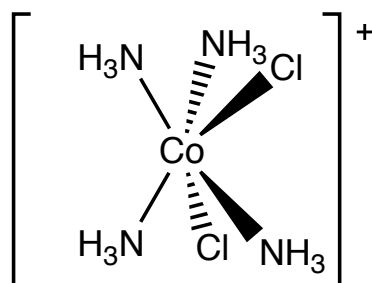
possible 6-coordinate models



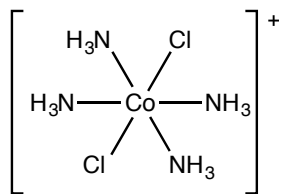
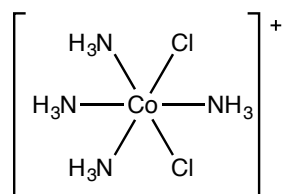
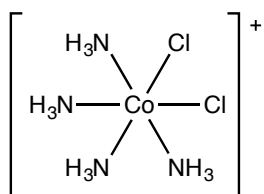
hexagonal

hexagonal
pyramidal

Octahedral

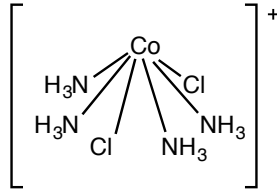
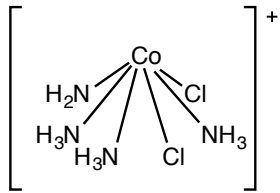
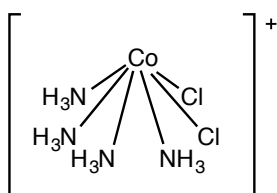
trigonal
antiprismatictrigonal
prismatic

possible 6-coordinate models

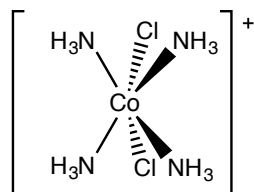
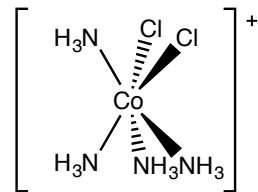
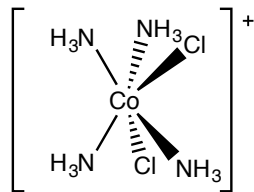


hexagonal

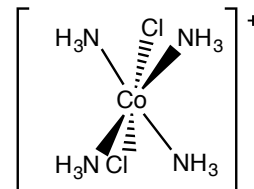
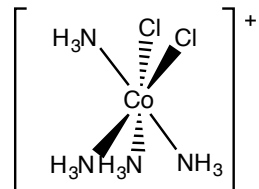
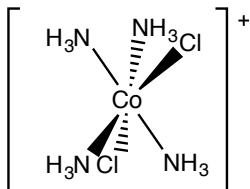
Models suggest 3 isomers, only 2 observed



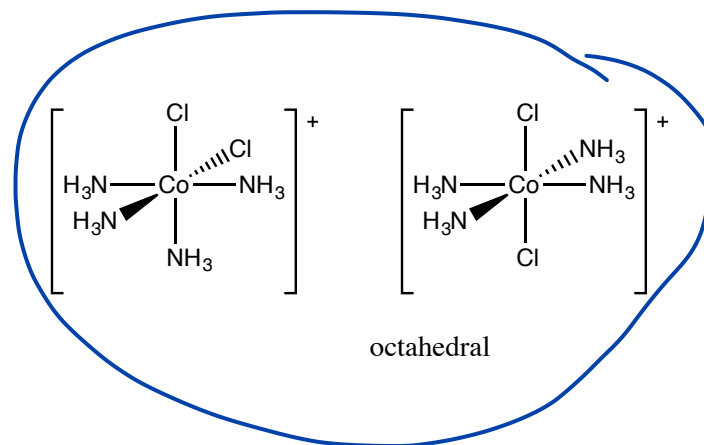
hexagonal pyramidal



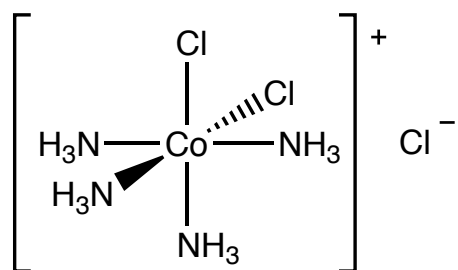
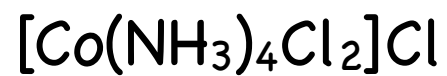
trigonal prismatic



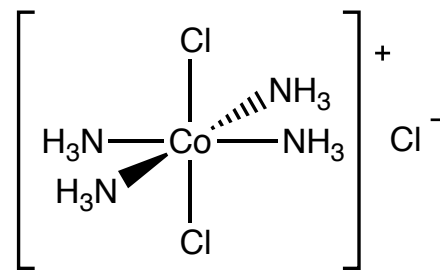
trigonal anti prismatic



octahedral



cis

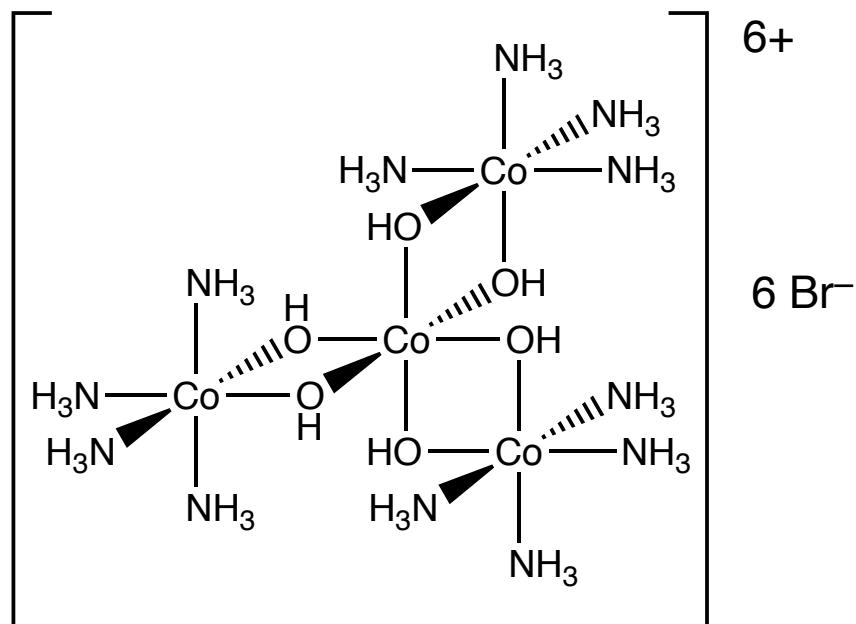


trans

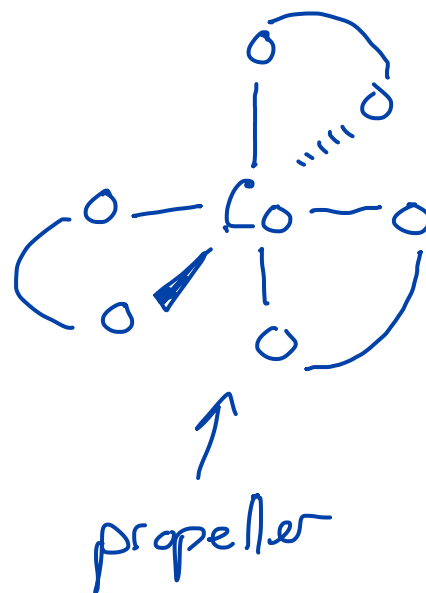
$$S_1 = \sigma$$

$$S_2 = i$$

$$S_3 \dots$$



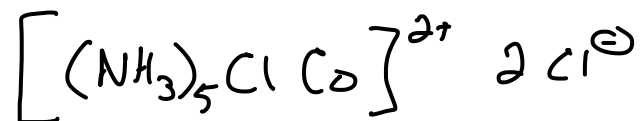
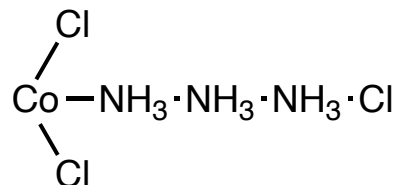
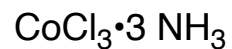
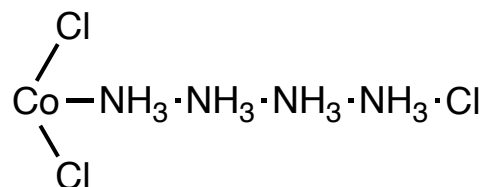
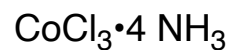
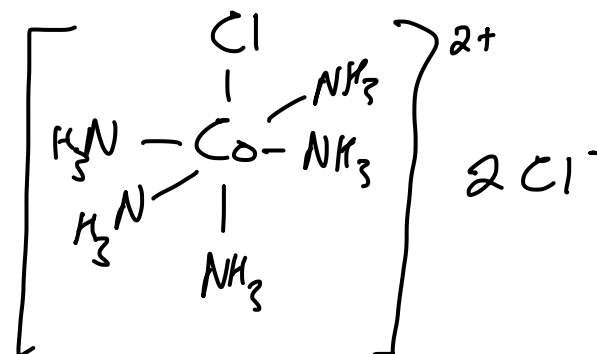
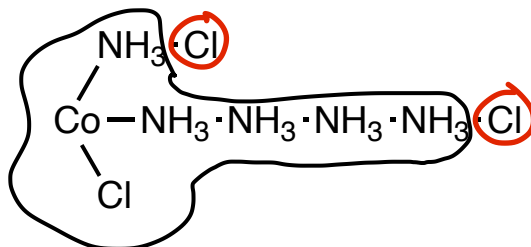
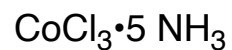
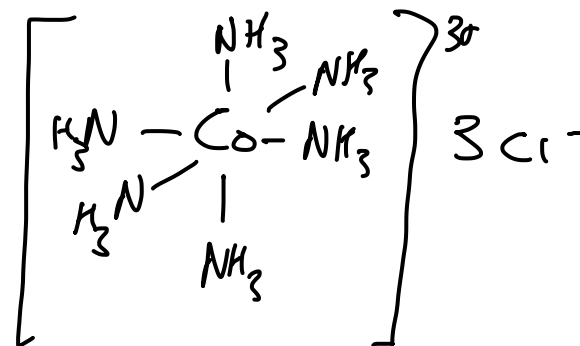
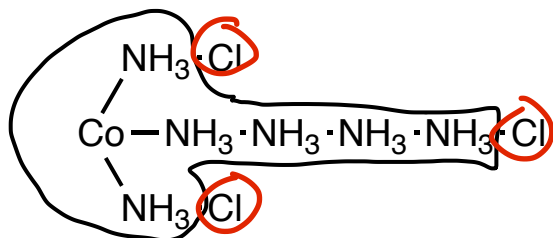
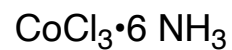
chiral



neither has
a plane
of
symmetry

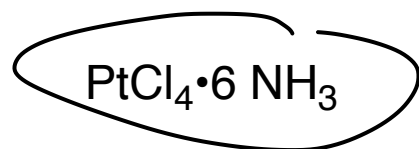
only the octahedral shape would create
a pair of enantiomers for this molecule

complex ion



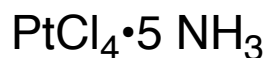
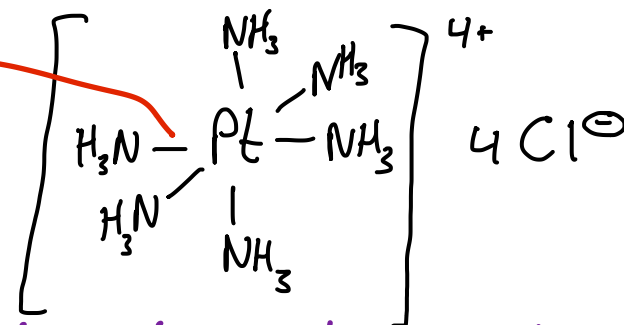
to precipitate 4 Cl^- 's the Cl^- 's must not be directly bonded to the octahedral metal (octahedral = 6 coordinate)

Cl^- ions precipitated by Ag^+



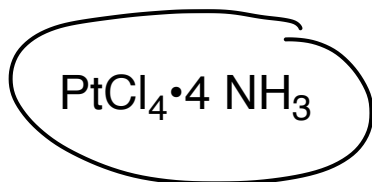
4

donative or coordinate covalent bond

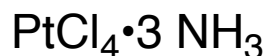
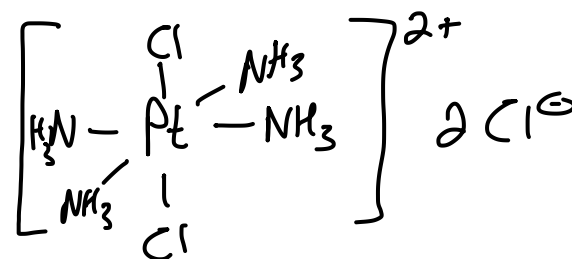


3

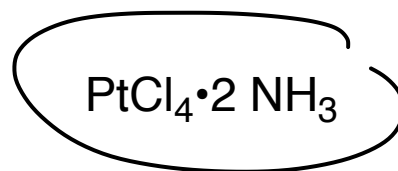
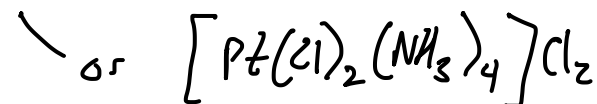
the N shares its e's but maintains ownership of both of them



2



1



0

