

1>2>3 % S orbital & En
1>2>3

(33) 3>2>1



(35) Carbamion $\text{SP}^2\oplus$ $\text{SP}^2\oplus$ $\text{SP}^2\oplus$

-I \downarrow O CH_3 -I \downarrow O CH_3 NO $_2$ \downarrow -I

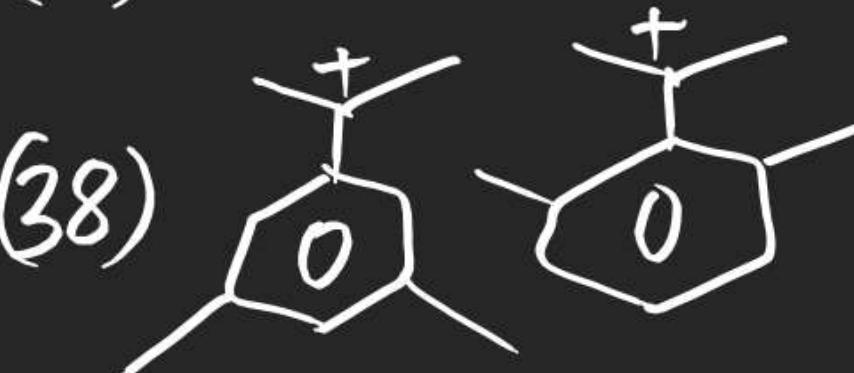
4>2>1>3

(36) $\text{SP}^2\oplus$ $\text{SP}^2\oplus$ $\text{SP}^2\oplus$ $\text{SP}^2\oplus$

+I \downarrow C H_3 -I \downarrow O CH_3 -I \downarrow O CH_3 NO $_2$ \downarrow -I

1>2>3>4

(37) Carbanion (4>3>2>1)



1>2 (SIR)

(38) Free Radical 1>2 (SIR)

(39) Carbanion 1>2 (SIR)

(40) Carbanion

(41) $\text{P}^{\oplus}\text{OCH}_3$ $\text{P}^{\oplus}\text{OCH}_3$ $\text{P}^{\oplus}\text{OCH}_3$ $\text{OCH}_3\text{P}^{\oplus}$

<+I <-I (1>3>2>4)

(42) Carbanion. (4>2>3>1)

(43)



(2 > 1) (Bredt's Rule)

(44)



(1 > 3 > 2)

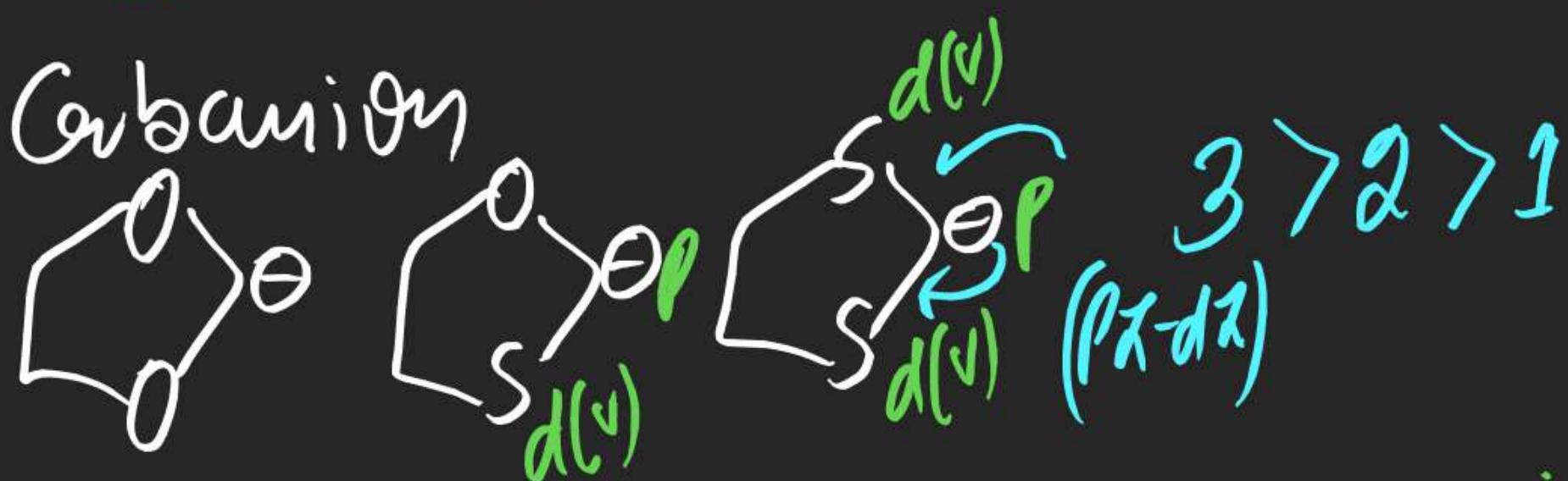
(45)



1 > 2 > 3
(2P-2P > 2P-3P > 3P-3P)

(46)

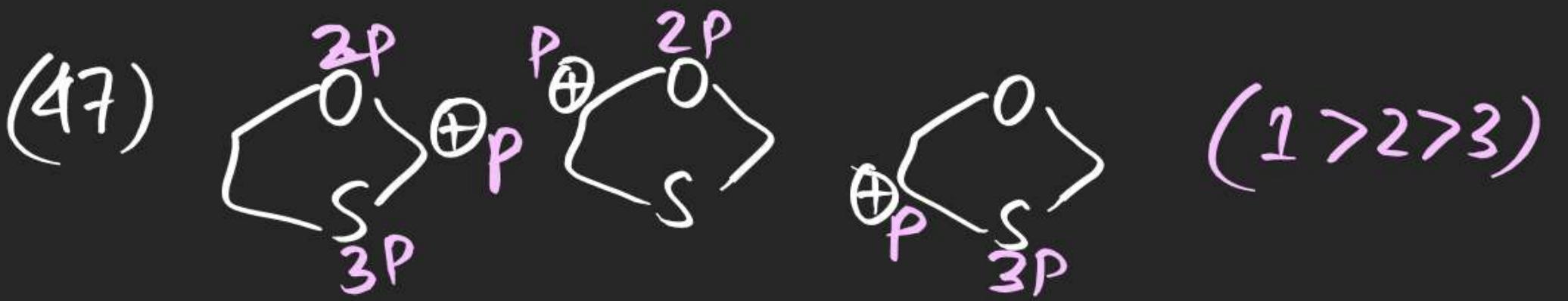
Carbamion



3 > 2 > 1

(Px-dz)

∴



(Stability
order)



Condition for Aromatic Compound :-

Compounds must be

- (a) Cyclic
- (b) planar (sp or sp^2)
- (c) Cyclic Conjugated
- (d) $(4n+2)\pi e\beta$ { $n=0, 1, 2, 3, \dots$ }
Nucleo β Rule Nucleo No. (2, 6, 10, 14, ...)

(#) Anti Aromatic Compounds:

All cyclic Compounds which are highly unstable than its open chain analogous system are known as Anti aromatic compound.



Condition for Anti Aromatic Compound

- (a) Ayclic ✓
- (b) Planar ✓
- (c) Conjugated ✓
- (d) $4n\pi$ electrons ($n=1, 2, 3, \dots$) [4, 8, 12, 16, ...]

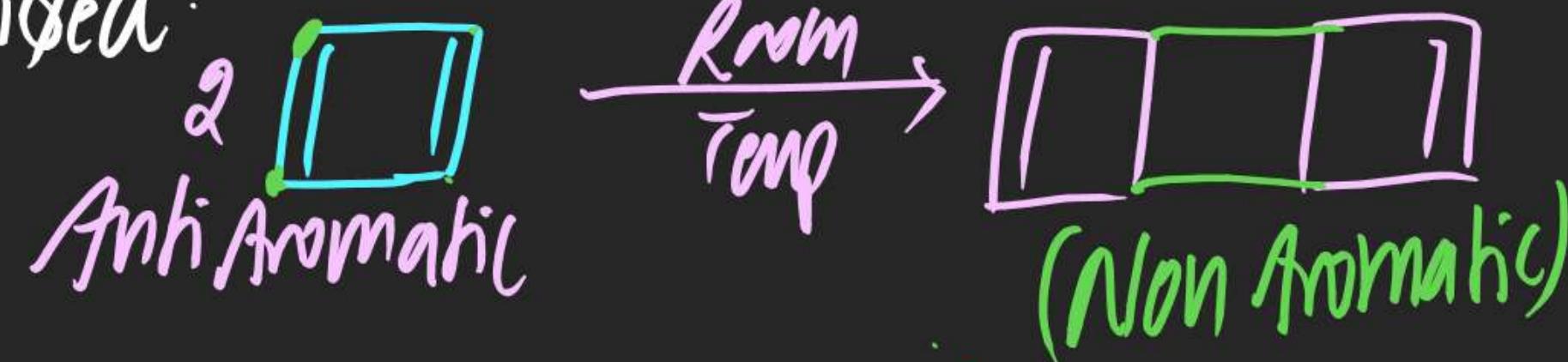
(#) Non Aromatic Compound:

⇒ Compounds which are neither Aromatic nor Anti Aromatic are known as Non Aromatic Compounds.

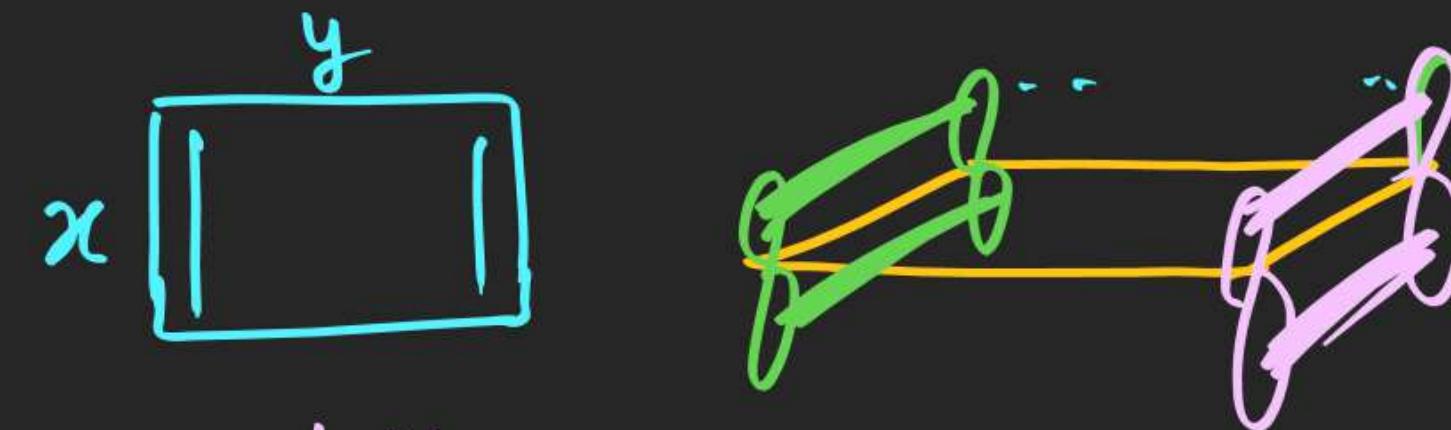
Note: (i) Stability order

~~Temp~~
~~(ii)~~ Aromatic Compound > Non Aromatic > Anti Aromatic
Cyclic Compounds containing more than 7 carbon atoms
are never Anti Aromatic

(iii) Anti Aromatic compound don't exist at Room Temperature in its pure form & gets dimerized.



In CycloButa-1,3-diene it is found that it contains two different C-C Bond lengths and it means AntiAromatic Compounds are not stabilised by Resonance.



$$x \neq y$$

(#) Quasi Aromatic Compound:

⇒ Aromatic Compounds having charge delocalisation are known as

Quasi Aromatic Compound

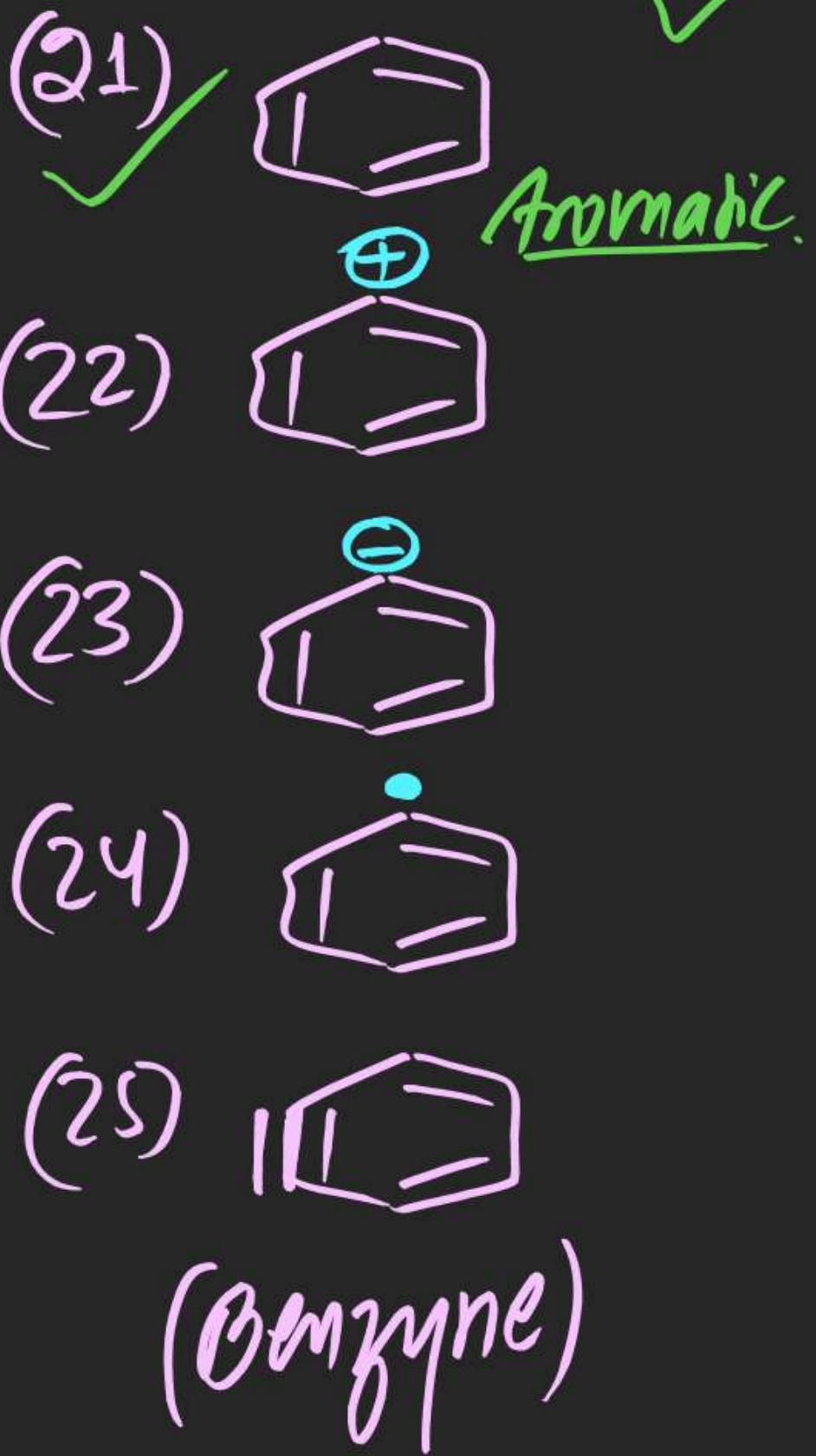


Cyclic Conjugated Planes

		Cyclic	Conjug. Planes	Hückel No.
(1)		✓	✗	✓ 2π
(2)		✓	✓	✓ 2π
(3)		✓	✓	✓ (2π)
(4)		✓	✓	4π
(5)		✓	✓	2π

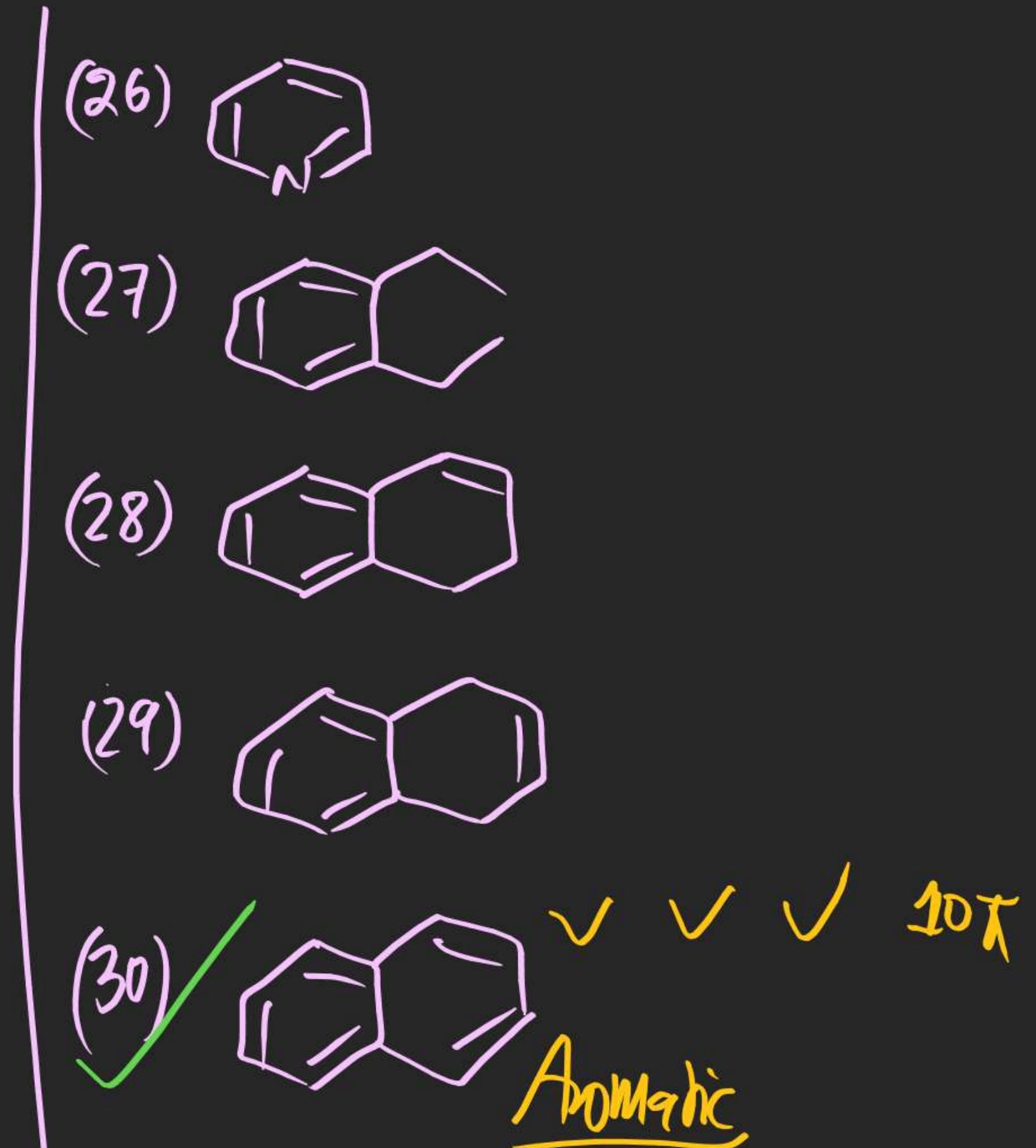
(6)		✓	✓	✓	4π Anti Aromatic
(7)		✓	✓	✓	2π Aromatic
(8)		✓	✓	✓	6π Aromatic
(9)		✓	✗	✗	2π Non Aromatic
(10)		✓	✗	✓	4π Non Aromatic





✓ ✓ ✓

6R



(31)



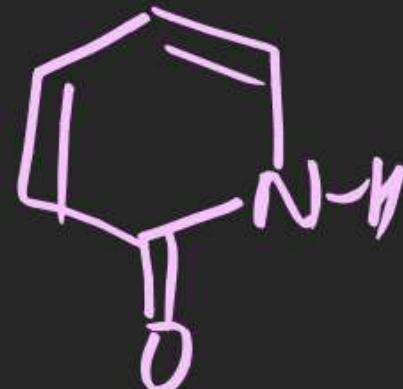
(32)



(33)



(34)



(35) Inorganic Benzene

or

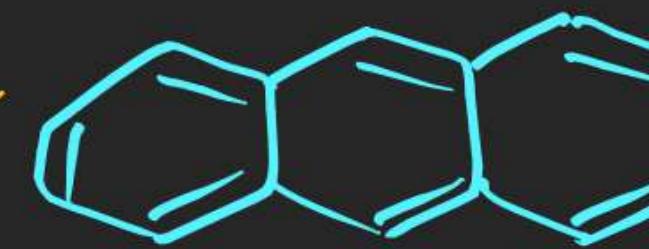
Borazine

($B_3N_3H_6$)

or

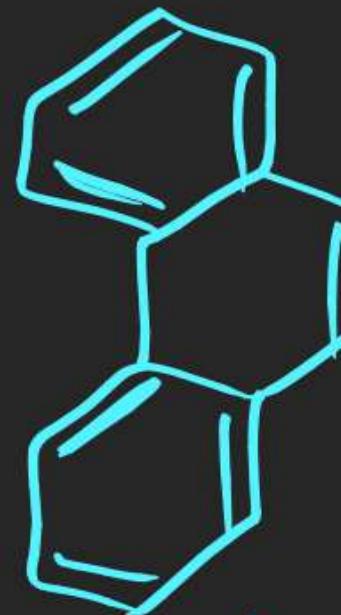
Borazole

(36)



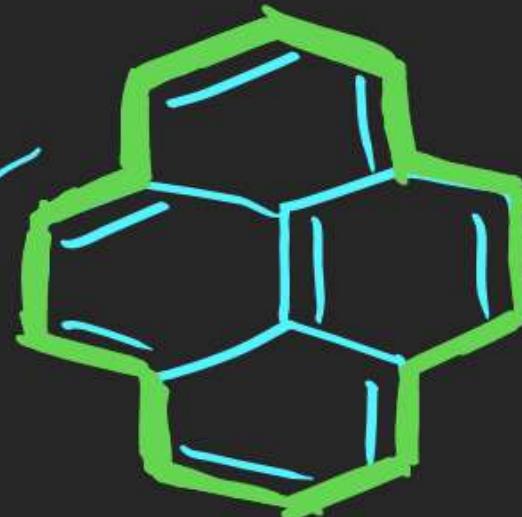
(Aromatic)

(37)



✓ ✓ ✓ 14

(38)

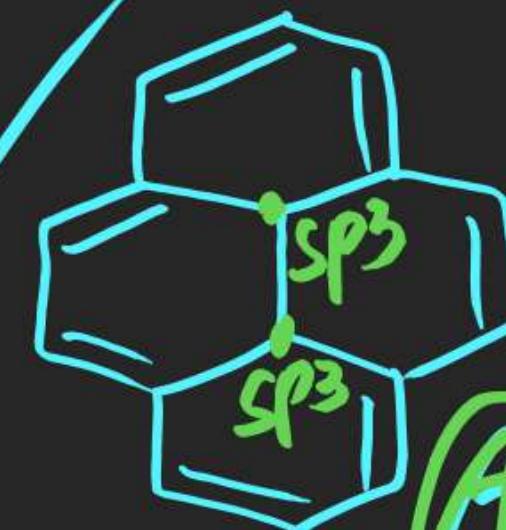


largest conjugated cyclic periphay.

(41)



(39)

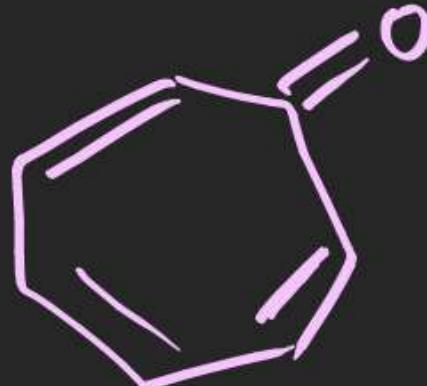


(Aromatic)

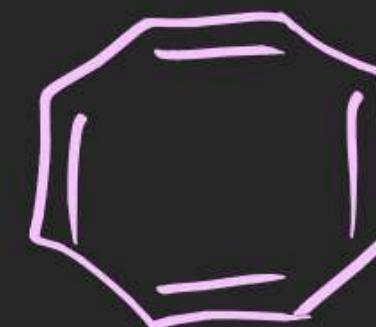
(40)



(42)

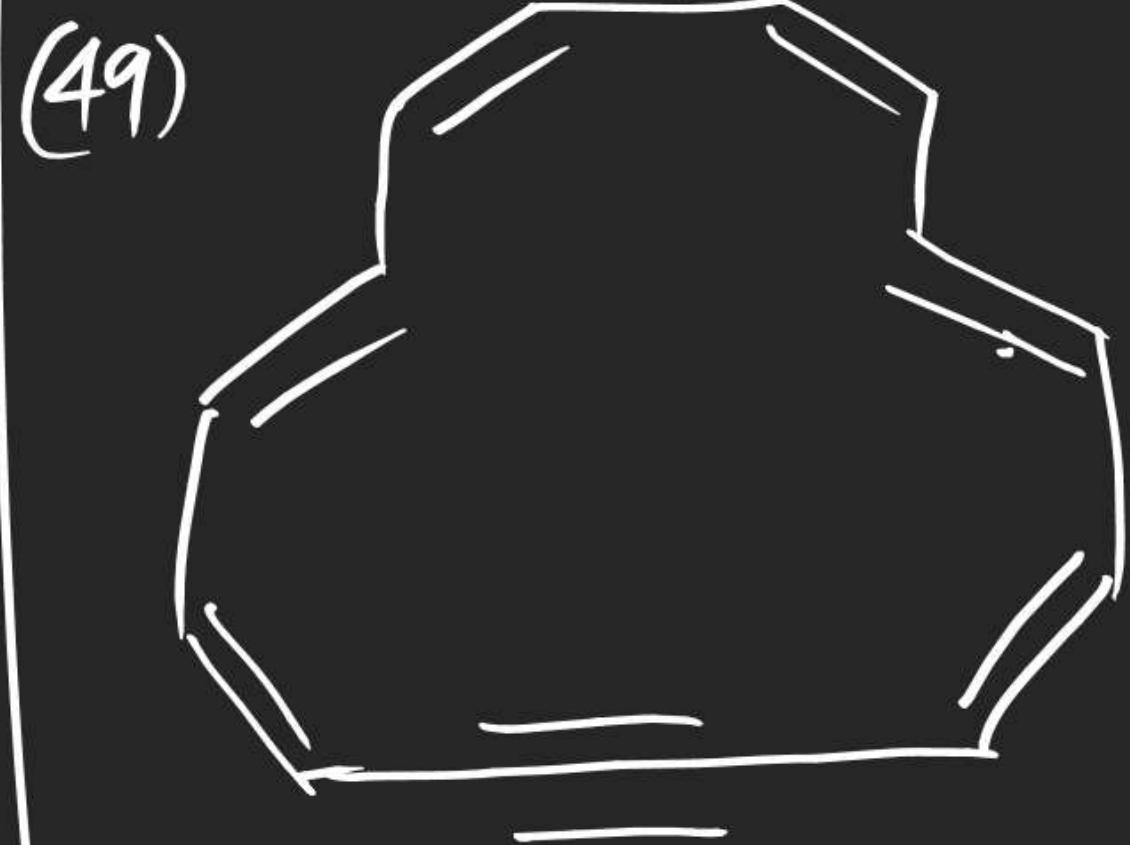
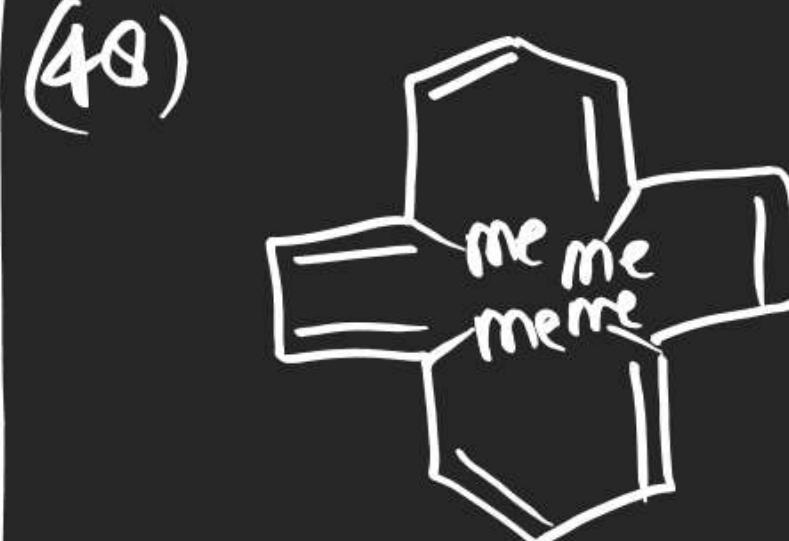


(43)



(44)





(50) Annulene-[16]

(51) Annulene-[18]

(52) Azulene

