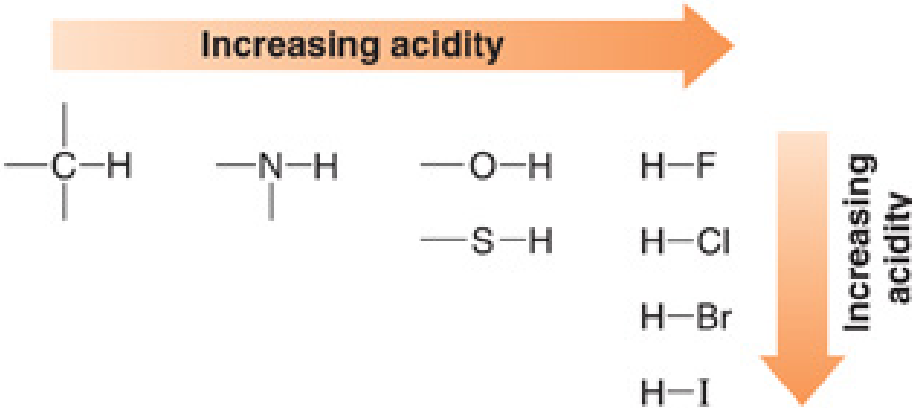



Fattori che determinano la forza acida

- **Qualsiasi parametro che stabilizzi una base coniugata A^- rende l'acido di partenza H-A più acido.**
- **Quattro fattori condizionano l'acidità di H-A. Essi sono:**
 - **Effetti dell'elemento A**
 - **Effetti induttivi**
 - **Effetti della risonanza**
 - **Effetti dell'ibridazione**
- **per confrontare l'acidità di due acidi confrontabili si segue sempre la stessa procedura:**
 - **Disegnare sempre le basi coniugate.**
 - **Determinare quale base coniugata sia più stabile.**
 - **Più è stabile la base coniugata, più forte è l'acido.**



Riassumendo:



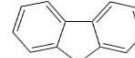
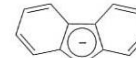
Factor	Example	
<p>1. Element effect: The acidity of H–A increases both left-to-right across a row and down a column of the periodic table.</p>	 <p style="text-align: center;">Increasing acidity →</p> <p style="text-align: center;"> $\begin{array}{cccc} \text{—C—H} & \text{—N—H} & \text{—O—H} & \text{H—F} \\ & & \text{—S—H} & \text{H—Cl} \\ & & & \text{H—Br} \\ & & & \text{H—I} \end{array}$ </p> <p style="text-align: right; vertical-align: middle;">↓ Increasing acidity</p>	
<p>2. Inductive effects: The acidity of H–A increases with the presence of electron-withdrawing groups in A.</p>	$\text{CH}_3\text{CH}_2\text{O—H}$	$\text{CF}_3\text{CH}_2\text{O—H}$ more acidic
<p>3. Resonance effects: The acidity of H–A increases when the conjugate base A^- is resonance stabilized.</p>	$\text{CH}_3\text{CH}_2\text{O—H}$	$\text{CH}_3\text{COO—H}$ more acidic
<p>4. Hybridization effects: The acidity of H–A increases as the percent s-character of A^- increases.</p>	CH_3CH_3 $\text{CH}_2=\text{CH}_2$ $\text{H—C}\equiv\text{C—H}$  <p style="text-align: center;">Increasing acidity →</p>	

Valori di pK_a di alcuni acidi

Acido	Base	pK_a
RNO_2^+	RNO_2	-12
ArNO_2^+	ArNO_2	-11
HClO_4	ClO_4^-	-10
HI	I^-	-10
RCNH^+	RCN	-10
$\text{R}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{H}$	$\text{R}-\overset{\text{O}}{\parallel}{\text{C}}-\text{H}$	-10
HBr	Br^-	-9
$\text{Ar}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{OR}$	$\text{Ar}-\overset{\text{O}}{\parallel}{\text{C}}-\text{OR}$	-7,4
HCl	Cl^-	-7
RSH_2^+	RSH	-7
$\text{Ar}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{OH}$	$\text{Ar}-\overset{\text{O}}{\parallel}{\text{C}}-\text{OH}$	-7
$\text{Ar}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{H}$	$\text{Ar}-\overset{\text{O}}{\parallel}{\text{C}}-\text{H}$	-7
$\text{R}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{R}$	$\text{R}-\overset{\text{O}}{\parallel}{\text{C}}-\text{R}$	-7
ArSO_3H	ArSO_3^-	-6,5
$\text{R}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{OR}$	$\text{R}-\overset{\text{O}}{\parallel}{\text{C}}-\text{OR}$	-6,5
ArOH_2^+	ArOH	-6,4
$\text{R}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{OH}$	$\text{R}-\overset{\text{O}}{\parallel}{\text{C}}-\text{OH}$	-6

Acido	Base	pK_a
$\text{Ar}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{R}$	$\text{Ar}-\overset{\text{O}}{\parallel}{\text{C}}-\text{R}$	-6
$\text{Ar}-\overset{\text{+}}{\text{O}}-\text{R}$	$\text{Ar}-\text{O}-\text{R}$	-6
$\text{CH}(\text{CN})_3$	$^-\text{C}(\text{CN})_3$	-5
Ar_3NH^+	Ar_3N	-5
$\text{H}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{H}$	$\text{H}-\overset{\text{O}}{\parallel}{\text{C}}-\text{H}$	-4
$\text{R}-\overset{\text{+}}{\text{O}}-\text{R}$	$\text{R}-\text{O}-\text{R}$	-3,5
$\text{R}_3\overset{\text{+}}{\text{C}}\text{OH}_2$	R_3COH	-2
$\text{R}_2\overset{\text{+}}{\text{C}}\text{OH}_2$	R_2CHOH	-2
$\text{RCH}_2\overset{\text{+}}{\text{O}}\text{H}_2$	RCH_2OH	-2
H_3O^+	H_2O	-1,74
$\text{Ar}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{NH}_2$	$\text{Ar}-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}_2$	-1,5
HNO_3	NO_3^-	-1,4
$\text{R}-\overset{\text{+}}{\text{C}}(\text{OH})-\text{NH}_2$	$\text{R}-\overset{\text{O}}{\parallel}{\text{C}}-\text{NH}_2$	-0,5
$\text{Ar}_2\overset{\text{+}}{\text{N}}\text{H}_2$	Ar_2NH	1
HSO_4^-	SO_4^{2-}	1,99
HF	F^-	3,17
HONO	NO_2^-	3,29
$\text{Ar}\overset{\text{+}}{\text{N}}\text{H}_3$	ArNH_2	3-5

Acido	Base	pK_a
$\text{Ar}\overset{+}{\text{N}}\text{R}_2\text{H}$	ArNR_2	3-5
RCOOH	RCOO^-	4-5
HCOCH_2CHO	$\text{HCO}\bar{\text{C}}\text{HCHO}$	5
H_2CO_3	HCO_3^-	6,35
H_2S	HS^-	7,00
ArSH	ArS^-	6-8
$\text{CH}_3\text{COCH}_2\text{COCH}_3$	$\text{CH}_3\text{CO}\bar{\text{C}}\text{HCOCH}_3$	9
HCN	CN^-	9,2
$^+\text{NH}_4$	NH_3	9,24
ArOH	ArO^-	8-11
RCH_2NO_2	$\text{R}\bar{\text{C}}\text{HNO}_2$	10
$\text{R}_3\overset{+}{\text{N}}\text{H}$	R_3N	10-11
$\text{R}\overset{+}{\text{N}}\text{H}_3$	RNH_2	10-11
HCO_3^-	CO_3^{2-}	10,33
RSH	RS^-	10-11
$\text{R}_2\overset{+}{\text{N}}\text{H}_2$	R_2NH	11
NCCH_2CN	$\text{NC}\bar{\text{C}}\text{HCN}$	11
$\text{CH}_3\text{COCH}_2\text{COOR}$	$\text{CH}_3\text{CO}\bar{\text{C}}\text{HCOOR}$	11
$\text{CH}_3\text{SO}_2\text{CH}_2\text{SO}_2\text{CH}_3$	$\text{CH}_3\text{SO}_2\bar{\text{C}}\text{HSO}_2\text{CH}_3$	12,5
$\text{EtOOCCH}_2\text{COOEt}$	$\text{EtOOC}\bar{\text{C}}\text{HCOOEt}$	13
CH_3OH	CH_3O^-	15,2
H_2O	OH^-	15,74
		16
RCH_2OH	RCH_2O^-	16

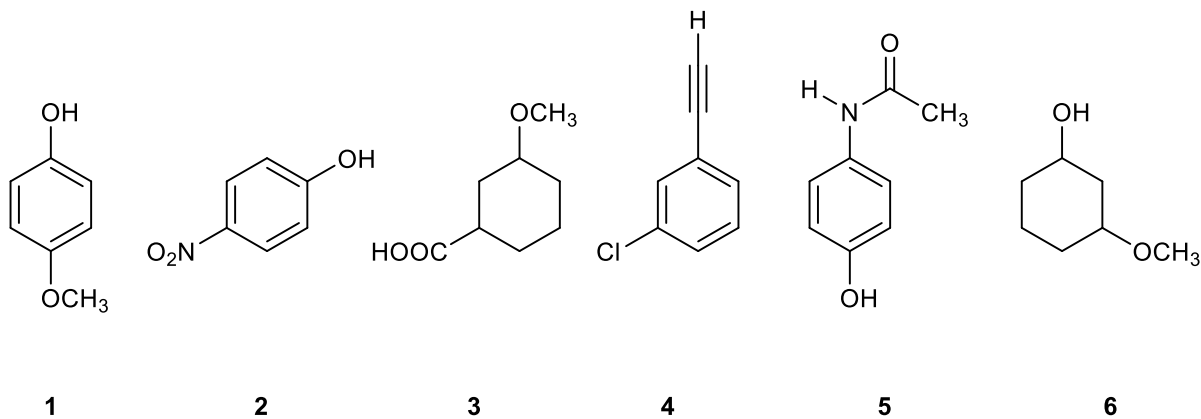
Acido	Base	pK_a
RCH_2CHO	$\text{R}\bar{\text{C}}\text{HCHO}$	16
R_2CHOH	R_2CHO^-	16,5
R_3COH	R_3CO^-	17
RCONH_2	RCONH^-	17
RCOCH_2R	$\text{RCO}\bar{\text{C}}\text{HR}$	19-20
		20
		23
ROOCCH_2R	$\text{ROOC}\bar{\text{C}}\text{HR}$	24,5
RCH_2CN	$\text{R}\bar{\text{C}}\text{HCN}$	25
$\text{HC}\equiv\text{CH}$	$\text{HC}\equiv\text{C}^-$	25
Ar_3CH	Ar_3C^-	31,5
Ar_2CH_2	Ar_2CH^-	33,5
H_2	H^-	35
NH_3	NH_2^-	38
PhCH_3	PhCH_2^-	40
$\text{CH}_2=\text{CHCH}_3$	$[\text{CH}_2^{\text{---}}\text{CH}^{\text{---}}\text{CH}_2]^-$	43
PhH	Ph^-	43
$\text{CH}_2=\text{CH}_2$	$\text{CH}_2=\text{CH}^-$	44
ciclo- C_3H_6	ciclo- C_3H_5^-	46
CH_4	CH_3^-	48
C_2H_6	C_2H_5^-	50
$(\text{CH}_3)_2\text{CH}_2$	$(\text{CH}_3)_2\text{CH}^-$	51

Gruppo	Struttura	+I	-I
alchilico	-R	*	
Alogeni	-X		*
idrossile	-OH		*
etereo (alcossi)	-OR		*
alcolato (alcossido)	-O ⁻	*	
carbossile	-COOH		*
ammidico	-CONH ₂		*
aldeidico	-COH		*
chetonico	-COR		*
amminico	-NH ₂		*
estereo	-COOR		*
nitro	-NO ₂		*
tiolato	-S ⁻	*	
tioetereo	-SR		*
ciano	-CN		*
solfidrilico (tiolico)	-SH		*
tioetereo	-SR		*

	Forza	Gruppo	Orientazione
Attivanti Elettron-donatori (EDG)	Forti	-O ⁻ -OH -NH ₂ -NHR -NR ₂	orto-para
	Medi	-OCH ₃ -OR -NHCOCH ₃ -NHCOR	orto-para
	Deboli	-CH ₃ -C ₂ H ₅ -R -C ₆ H ₅ -CH=CR ₂	orto-para
Riferimento	Neutro	-H	
Disattivanti Elettron-attrattori (EWG)	deboli	-F -Cl -Br -I	orto-para
	Medi	-COH, -COR -COOH, COOR -COX, -CONH ₂ -SO ₃ H -CN	meta
	Forti	-CF ₃ , CCl ₃ ,.. -NH ₃ ⁺ -NR ₃ ⁺ -NO ₂	meta

ESERCITAZIONI: Scale acidità e basicità_1

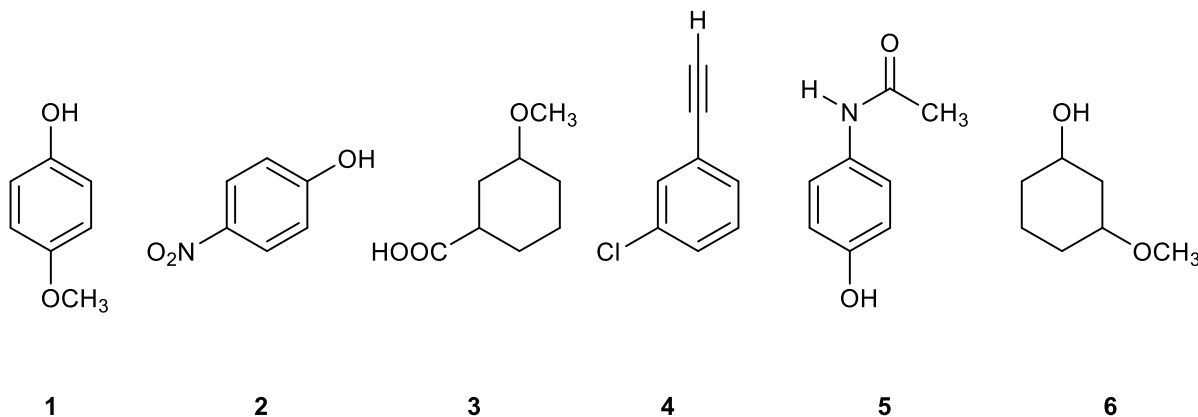
1. Indicare il corretto ordine di acidità crescente



- A** 4<6<3<1<5<2
B 6<5<2<1<4<3
C 4<1<5<2<6<3
D 2<5<1<4<3<6
E 4<6<1<5<2<3

ESERCITAZIONI: Scale acidità e basicità_1

1. Indicare il corretto ordine di acidità crescente



A 4<6<3<1<5<2

B 6<5<2<1<4<3

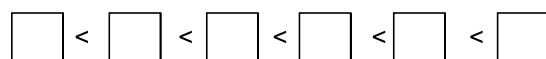
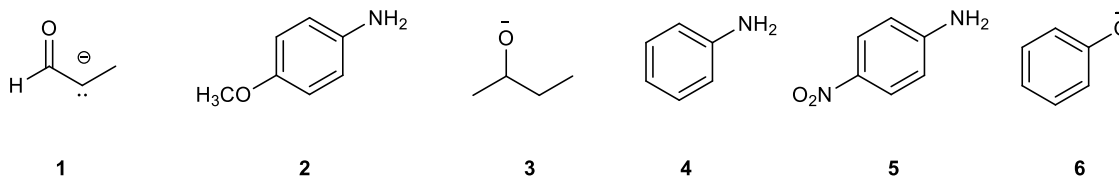
C 4<1<5<2<6<3

D 2<5<1<4<3<6

X 4<6<1<5<2<3

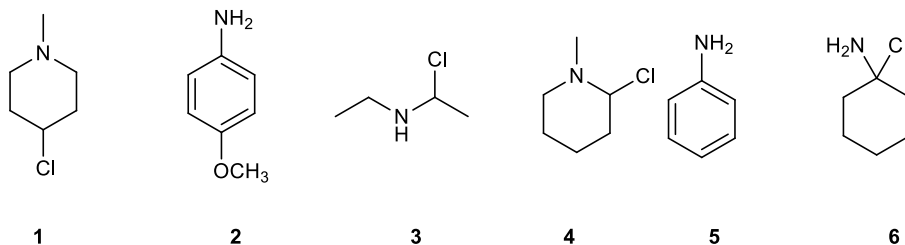
ESERCITAZIONI: Scale acidità e basicità_2

1.



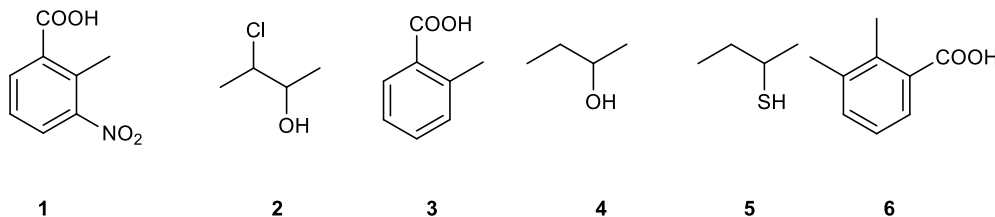
basicità crescente

2. Indicare il corretto ordine di basicità



- A** 5>2>3>6>4>1
- B** 4>1>3>6>2>5
- C** 4>1>6>3>5>2
- D** 1>4>3>6>2>5
- E** 1>4>6>2>5>3

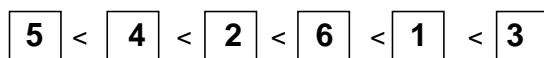
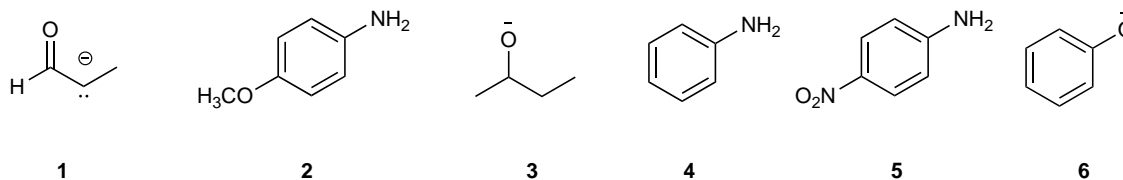
3. Indicare il corretto ordine di acidità crescente



- A** 1>6>3>5>2>4
- B** 3>1>6>5>2>4
- C** 1>3>6>5>2>4
- D** 1>3>6>2>4>5
- E** 1>3>6>5>4>2

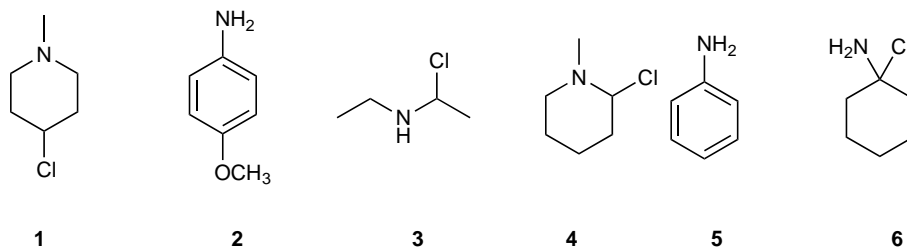
ESERCITAZIONI: Scale acidità e basicità_Soluzioni_2

1.



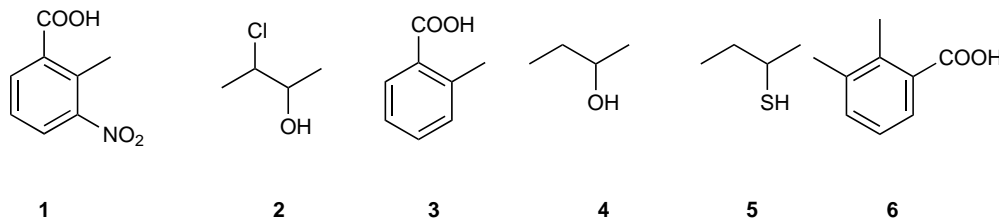
basicità crescente

2. Indicare il corretto ordine di basicità



- A 5>2>3>6>4>1
- B 4>1>3>6>2>5
- C 4>1>6>3>5>2
- ✖ D 1>4>3>6>2>5
- E 1>4>6>2>5>3

3. Indicare il corretto ordine di acidità crescente



- A 1>6>3>5>2>4
- B 3>1>6>5>2>4
- ✖ C 1>3>6>5>2>4
- D 1>3>6>2>4>5
- E 1>3>6>5>4>2