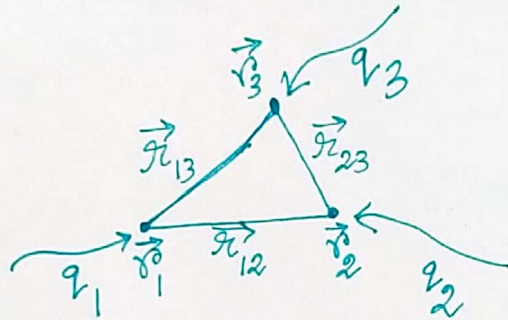


$$\vec{\nabla} \cdot (f \vec{A}) = f (\vec{\nabla} \cdot \vec{A}) + \vec{A} \cdot (\vec{\nabla} f)$$

$$\Rightarrow \int \vec{\nabla} \cdot (f \vec{A}) dV = \int f (\vec{\nabla} \cdot \vec{A}) dV + \int \vec{A} \cdot (\vec{\nabla} f) dV$$

$$\Rightarrow \boxed{\int f (\vec{\nabla} \cdot \vec{A}) dV = - \int \vec{A} \cdot (\vec{\nabla} f) dV + \oint (f \vec{A}) \cdot d\vec{S}}$$

Energy of a point charge distribution:



Work done to bring q_1 to $\vec{r}_1 = 0$

[Because, there was no electric field initially]

Once, q_1 is in place, it'll create an electrostatic field.

So, now when you try to bring q_2 to \vec{r}_2 , you need to do work against this field. Similarly when you are bringing q_3 to \vec{r}_3 , you need to work against the combined field created by q_1 and q_2 .

$$\text{Work done to bring } q_2 \text{ to } \vec{r}_2 = W_2 = \frac{1}{4\pi\epsilon_0} q_2 \left(\frac{q_1}{r_{12}} \right)$$

$$\text{" " " " } q_3 \text{ to } \vec{r}_3 = W_3 = \frac{1}{4\pi\epsilon_0} q_3 \left(\frac{q_1}{r_{13}} + \frac{q_2}{r_{23}} \right)$$

$$\text{" " " " } q_4 \text{ to } \vec{r}_4 = W_4 = \frac{1}{4\pi\epsilon_0} q_4 \left(\frac{q_1}{r_{14}} + \frac{q_2}{r_{24}} + \frac{q_3}{r_{34}} \right)$$

$$\Rightarrow W_4 = \frac{1}{8\pi\epsilon_0} \left[\frac{q_4 q_1}{r_{14}} + \frac{q_4 q_2}{r_{24}} + \frac{q_4 q_3}{r_{34}} + \frac{q_1 q_4}{r_{14}} + \frac{q_2 q_4}{r_{24}} + \frac{q_3 q_4}{r_{34}} \right]$$

In general, $W_i = \frac{1}{8\pi\epsilon_0} q_i \left(\sum_{j \neq i} q_j / r_{ij} \right)$

Total work done to assemble n charges $= W = \frac{1}{8\pi\epsilon_0} \sum_{i=1}^n \sum_{\substack{j=1 \\ (i \neq j)}}^n \frac{q_i q_j}{r_{ij}}$

①

————— (1)

$$W = \frac{1}{2} \sum_{i=1}^n q_i \left(\sum_{\substack{j=1 \\ j \neq i}}^n \frac{1}{4\pi\epsilon_0} \frac{q_j}{r_{ij}} \right)$$

$$\Rightarrow \boxed{W = \frac{1}{2} \sum_{i=1}^n q_i V(r_i)} \quad \text{--- (2)}$$

$V(r_i)$ = potential at r_i due to all charges other than q_i .

Energy of a continuous charge distribution:

$$W = \frac{1}{2} \int_V \rho V d\tau \quad \text{--- (3)}$$

$$\rho = \epsilon_0 (\nabla \cdot \vec{E}). \text{ From (3),}$$

$$\begin{aligned} W &= \frac{\epsilon_0}{2} \int_V (\nabla \cdot \vec{E}) V d\tau \\ &= \frac{\epsilon_0}{2} \left[- \int_V \vec{E} \cdot (\nabla V) d\tau + \oint_S V \vec{E} \cdot d\vec{S} \right] \end{aligned}$$

[We have used: $\int_V f(\nabla \cdot \vec{A}) d\tau = - \int_V \vec{A} \cdot (\nabla f) d\tau + \oint_S f \vec{A} \cdot d\vec{S}]$

$$\therefore W = \frac{\epsilon_0}{2} \left[\int_V E^2 d\tau + \oint_S V \vec{E} \cdot d\vec{S} \right] \quad \text{--- (4)}$$

But what volume and surface were we talking about?
The volume V must enclose all the charges.
However, from eq.(3), if we include a larger volume, $\rho = 0$ ensures that there's no contribution to W beyond V .

Q. Does W diverge if we increase the volume?
 E^2 is positive. So, the first term in eq.(4) increases.
But the second term must decrease in order to leave the sum intact.

But how to understand this?

$$V \sim \frac{1}{r^0} \text{ (at least)}, E \sim \frac{1}{r^{0.2}} \text{ (at least)}$$

$$\text{So, } EV \sim \frac{1}{r^{0.3}} \text{ but, } S \sim r^{0.2}$$

$$\therefore EV ds \sim \frac{1}{r^0}$$

$$\text{As } r \rightarrow \infty, \oint \vec{E} \cdot d\vec{S} \rightarrow 0$$

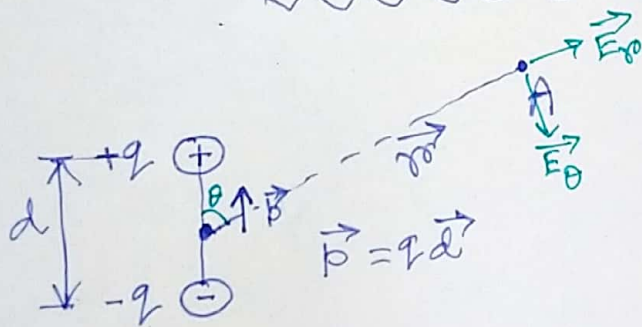
$$\therefore W = \frac{\epsilon_0}{2} \int_{\text{all space}} E^2 dV$$

For any physical field, $E \rightarrow 0$ as $r \rightarrow \infty$.

$$\therefore W = \frac{\epsilon_0}{2} \int E^2 dV \text{ doesn't diverge!}$$

$$\text{Energy per unit volume} = \frac{1}{2} \epsilon_0 E^2$$

Electrostatic field inside a material



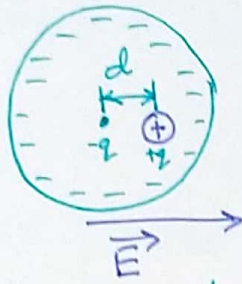
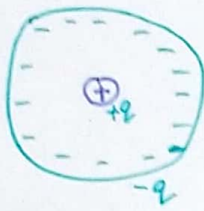
$$V(r, \theta) = \frac{1}{4\pi\epsilon_0} \frac{p \cos\theta}{r^2} = \left(\frac{\vec{p} \cdot \hat{r}}{4\pi\epsilon_0 r^2} \right) \quad (5)$$

$$\begin{aligned} \vec{E} &= -\vec{\nabla} V(r, \theta) = -\left(\hat{r} \frac{\partial V}{\partial r} + \hat{\theta} \frac{1}{r} \frac{\partial V}{\partial \theta} \right) \\ &= \hat{r} \frac{2p \cos\theta}{4\pi\epsilon_0 r^3} + \hat{\theta} \frac{p \sin\theta}{4\pi\epsilon_0 r^3} \end{aligned}$$

$$E_r = \frac{2p \cos\theta}{4\pi\epsilon_0 r^3}, \quad E_\theta = \frac{p \sin\theta}{4\pi\epsilon_0 r^3}$$

(3)

Induced dipole:



$$\vec{p} \propto \vec{E}$$

$$\Rightarrow \vec{p} = \alpha \vec{E}$$

(α is called polarizability)

This is what happens to an atom when you place it in an E-field.

Same thing happens to molecules as well, but the situation may be slightly complicated. Consider, CO_2 :



If \vec{E} is along the axis (represented by E^{\parallel}) polarizability is more than if we had applied \vec{E} perpendicular to the axis! (i.e. $\alpha^{\parallel} > \alpha^{\perp}$)

$$\vec{p} = \alpha^{\parallel} \vec{E}^{\parallel} + \alpha^{\perp} \vec{E}^{\perp}$$

In general,

$$\begin{bmatrix} p_x \\ p_y \\ p_z \end{bmatrix} = \begin{bmatrix} \alpha_{xx} & \alpha_{xy} & \alpha_{xz} \\ \alpha_{yx} & \alpha_{yy} & \alpha_{yz} \\ \alpha_{zx} & \alpha_{zy} & \alpha_{zz} \end{bmatrix} \begin{bmatrix} E_x \\ E_y \\ E_z \end{bmatrix}$$

$$\begin{aligned} \therefore p_x &= \alpha_{xx} E_x + \alpha_{xy} E_y + \alpha_{xz} E_z \\ p_y &= \alpha_{yx} E_x + \alpha_{yy} E_y + \alpha_{yz} E_z \\ p_z &= \alpha_{zx} E_x + \alpha_{zy} E_y + \alpha_{zz} E_z \end{aligned}$$

↖ polarizability "matrix" (tensor, to be formal)

$$\vec{p} = [\alpha] \vec{E}$$

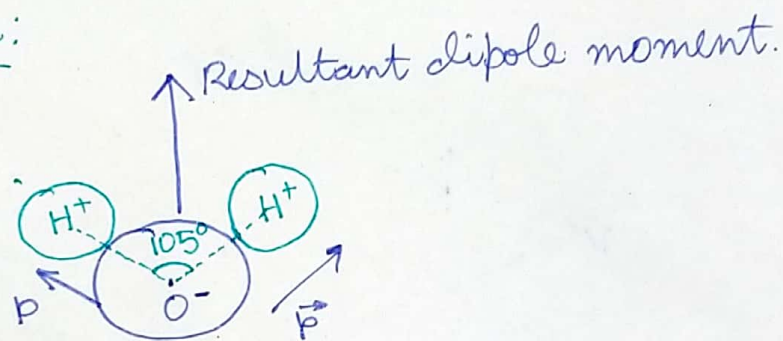
In general, \vec{p} and \vec{E} are not along the same direction.
 \vec{p} and \vec{E} are ~~only~~ along the same direction only if:

$$\begin{bmatrix} p_x \\ p_y \\ p_z \end{bmatrix} = \begin{bmatrix} \alpha & 0 & 0 \\ 0 & \alpha & 0 \\ 0 & 0 & \alpha \end{bmatrix} \begin{bmatrix} E_x \\ E_y \\ E_z \end{bmatrix} \leftarrow \text{Isotropic}$$

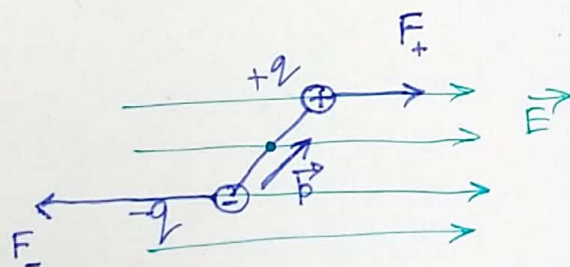
i.e. $p_x = \alpha E_x$, $p_y = \alpha E_y$, $p_z = \alpha E_z$

$$\therefore \vec{p} = \hat{x} p_x + \hat{y} p_y + \hat{z} p_z = \alpha [\hat{x} E_x + \hat{y} E_y + \hat{z} E_z] \\ = \alpha \vec{E}$$

Polar molecule:



What happens when such tiny dipoles are placed in an E-field?



$$\begin{aligned} \text{Torque}(\tau) &= \mathbf{r}_+ \times \vec{F}_+ + (\vec{r}_- \times \vec{F}_-) \\ &= \left(\frac{\vec{d}}{2} \times q\vec{E} \right) + \left\{ \left(-\frac{\vec{d}}{2} \right) \times (-q\vec{E}) \right\} \\ &= (q\vec{d} \times \vec{E}) = (\vec{p} \times \vec{E}) \end{aligned}$$

This torque'll tend to ~~align~~ align the dipole along the direction of \vec{E} .