Field of a shell via direct integration, for the field

For a point charge, $V=-\int \frac{1}{41160} \cdot \frac{9}{70248} \hat{r} \cdot dr \hat{r}$

$$= - kq \int_{\alpha}^{r} r^{-2-8} dr$$

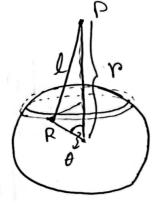
$$= - kq \cdot \frac{r^{-1-8}}{r^{-1-8}} \int_{\alpha}^{r}$$

$$= \frac{\kappa_{9}}{1+\delta} \cdot \frac{1}{\gamma^{1+\delta}} \Big|_{\infty}^{6 \gamma}$$

-.
$$V(r) = Kq \cdot \frac{1}{(4+6) r^{3+6}}$$

Now. let's calculate the potential due to a spherical non-conducting shell at a distace or from the center.

The infinitesimal area of the ring can be written as, dA = 211Rsind (Rdd)



Now,
$$dV(r) = K \frac{dq}{(4+8)q^{1+8}} = K \frac{\sigma 2\pi R^2 \sin\theta d\theta}{(4+8)q^{1+8}}$$

If
$$8=0$$
, then $\phi(r) = \frac{kB}{2rR} \cdot 2R = \frac{kB}{r} \cdot r R$
 $\phi(r) = \frac{kB}{2rR} \cdot 2r = \frac{kB}{R} \cdot r R R$

(b) The potential at on the swiface of the shell?

ignoring (1-82) torm,

ignoring
$$(1-8^2)$$
 torm,
$$\phi_{a_{1}a} = \frac{KB_a}{2a \cdot a} \cdot \left[f(a+a) - f(a-a) \right]$$

Similarly the potential of the soft shell with radius b

$$\phi_{a,b} = \frac{kB_b}{2ab} \left[f(a+b) - f(a-b) \right]$$

$$\frac{1}{2a} = \frac{kB_a}{2a^2} f(ba) + \frac{kB_b}{2ab} \left[f(a+b) - f(a-b) \right] 0$$

Similarly,
$$\phi_b = \frac{k\theta_b}{2b^2} f(2b) + \frac{k\theta_a}{2ab} \left[f(a+b) - f(a-b) \right]$$

1 For equipotential,

$$\phi_a = \phi_b = \phi$$

Multiplying (i) by a [flaty-flaty] and (ii) by bf(2a)

and subtracting -

$$\therefore \mathcal{G}_{b} = \frac{2b\phi}{k} \cdot \frac{bf(2a) - a \left[f(a+b) - f(a+b)\right]^{2}}{f(2a) f(2b) - \left[f(a+b) - f(a-b)\right]^{2}}$$

$$\text{If } \delta = 0, \text{ then } f(x) = x^{1-\delta} = x$$

$$\therefore \mathcal{G}_{b} = \frac{2b\phi}{k} \cdot \frac{2ab - a \times 2b}{-a \times 2b}$$

$$\therefore \mathcal{G}_{b} = 0$$

Work and energy in electroptation

Say, you have a configuration of source charges, and you want to move a test charge of from a to to to. We will have apply a force which is equal to the electrostatic force, but in apposite direction. In this way, you apply the minimum force. You could apply morre force than the electrostatic force, but that would increase the relocity and kinetic energy. We are only interested in the minimum force needed.

$$W = \int_{a}^{b} \vec{F} \cdot d\vec{s} = \int_{a}^{b} - B\vec{F} \cdot d\vec{s} = -9 \int_{a}^{b} \vec{F} \cdot d\vec{s}$$

$$= 9 \left[v(b) - v(b) \right]$$

$$V(b) - V(a) = \frac{W}{9}$$

Now, if you want to bring Q from very face away to a point P^3 , then, the work you must do is $W = Q[V(\vec{r}) - V(\omega)]$

: W= QV(F)

This is the work you will need, which is basically of times the potential. This can also be quantified as the potential energy, as we will get this much work/ energy back when the system is dismandled.

Energy of a point charge distribution

We want to calculate the work needed to assemble an entire collection of point charger. For the first charge q_1 there is no field and we do not need any work to to assemble it. For bringing q_2 , at a position \vec{E} , where q_1 is at \vec{r} is given b we need,

For bringing a third charge (keeping 9, and 8 fixed), the

$$W_{3} = q_{3} \bigvee_{3} (\vec{\xi}) + q_{3} \bigvee_{3} (\vec{r}_{3})$$

$$= q_{3} \bigvee_{3} \frac{q_{1}}{r_{13}} + q_{3} \frac{kq_{2}}{r_{23}}$$

For .. Total work = $K \frac{q_1 q_2}{Y_{12}} + K \frac{q_2 q_3}{Y_{23}} + K \frac{q_3 q_3}{Y_{31}}$

For a fourth charge Ta, we will have in total AC, terms in our total work. In general,

 $W = K \sum_{i=1}^{N} \sum_{j > i} \frac{q_i l_j}{r_{ij}}$

is making owne we are not counting. twice.

If we count all the pains twice, then -

$$W = \frac{1}{8\pi\epsilon} \cdot \sum_{i=1}^{n} \frac{n}{j=1} \frac{q_i q_{\dot{i}}}{\gamma_{i\dot{j}}^2}$$

$$\longrightarrow W = \frac{1}{2} \sum_{i=1}^{N} q_{i} \left[\sum_{j=1}^{N} \frac{1}{4\pi\epsilon} \frac{q_{j}}{\gamma_{i}} \right]$$

$$V(\vec{r}_i) = \frac{1}{2} \sum_{i=1}^{N} q_i V_i(\vec{r}_i)$$
 with $V(\vec{r}_i) = \sum_{\hat{\sigma}=1}^{N} \frac{\kappa q_{\hat{\sigma}}}{r_{\hat{\sigma}\hat{\sigma}}}$

The term in the parantheses is the potential at point i? (where 9; is located) due to all other charges. This is the energy stored in a configuration.

The energy of a continuous charge distribution

We can write, $\overrightarrow{\nabla} \cdot \overrightarrow{E} = + \frac{S}{\epsilon}$.

Now, in vector calculus, there is a vector identity, $\nabla \cdot (\nabla f) = f \nabla \cdot \vec{p} + \nabla f \cdot \vec{p}$ where \vec{p} is a vector field and f is a scalar function.

$$W = \frac{C}{2} \int \left(\overrightarrow{\nabla} \cdot (\overrightarrow{E} V) - \overrightarrow{\nabla} \overrightarrow{V} \cdot \overrightarrow{E} \right) d\mathcal{E}$$

$$= \frac{C}{2} \left[- \int_{-\overrightarrow{E}} \overrightarrow{E} d\mathcal{E} + \int_{-\overrightarrow{E}} \overrightarrow{\nabla} \cdot (\overrightarrow{V} \overrightarrow{E}) d\mathcal{E} \right]$$

$$= \frac{C}{2} \left[\int_{-\overrightarrow{E}} E^{2} d\mathcal{E} + \int_{-\overrightarrow{E}} \overrightarrow{\nabla} \cdot (\overrightarrow{V} \overrightarrow{E}) d\mathcal{E} \right]$$

Now, consider the volume where we are integrating. Initially we started with, $W = \frac{1}{2} | S \vee d ?$. So, the integration was over the space where the charges are located. But, we could expand the boundary as large as we want, up to infinity, because in that region, S = 0 and it contributes nothing to the total integral. Now, as we increase the volume, the volume integral increases, while the surface integral must decrease to keep W constant. The volume integral increases since we are integrating over the whole states of the volume. So, the integral increases if we pick a larger volume. So, the integral increases if we pick a larger volume. However, the surface integral is only computed on the boundary which grown like Φ r? where

Ex $\frac{1}{p_2}$ and $Vx \frac{1}{p}$. So, roughly speaking, the surface integral drops of as $\frac{1}{p}$ and if the volume is all space, then the surface integral goes to zero.

$$W = \frac{\mathcal{E}_{0}}{2} \underbrace{\mathbb{R}^{2}}_{\text{Mole}} \underbrace{\mathbb{R}^{2}}_{\text{Apace}}$$

$$W = \frac{\mathcal{E}_{0}}{2} \underbrace{\mathbb{R}^{2}}_{\text{Apace}} \underbrace{\mathbb{R}^{2}}_{\text{Apace}}$$

$$= \frac{2^{2}}{4\pi \mathcal{E}_{0}} \underbrace{\mathbb{R}^{2}}_{\text{Apace}} \underbrace{\mathbb{R}^{2}}_{\text{Apace}}$$

$$= \frac{4\pi q^{2}}{16\pi^{2}} \underbrace{\mathbb{R}^{2}}_{\text{Apace}} \underbrace{\mathbb{R}^{2}}_{\text{Apace}}$$

If R2 increases, the integral also increases. One could argue then that the electrophatic energy is stored in the electric field with energy density $U'=\frac{E_0}{Z}$ E^2 .

Now, there is a particular inconsistency between two equations—

First equation tells you that energy can be both positive and negative. Say, for example, the energy of two equal and opposite charges at a ditance

To apart in given by,
$$W = \frac{1}{2} \left[\frac{1}{4} V(r_1) + \frac{1}{2} V(r_2) \right]^{\frac{2}{r_1}} \frac{1}{r_2}$$

$$= \frac{1}{2} \left[\frac{1}{4 \text{Re}} \cdot \frac{-q}{r} + (-q) \cdot \frac{q}{4 \text{Re} r} \right]$$

$$= -\frac{1}{4 \text{Re}} \cdot \frac{q^2}{r^2}$$

However, second equation says energy is always positive Actually, they both are contribution of energy to make we didn't take the contribution of energy to make the point charges. We merely calculated the work done for moving -q to a distance in from 9 from infinity, where, the charges were given readymade. But, equation to second equation gives you energy of the whole configuration. But, there still is a problem with the second equation. If we calculate the total energy of a point charge, then,

But this doesn't make any sense, right? Where in the problem? The problem actually creet into one flace. In the equation, $W = \frac{1}{2} \sum_{i=1}^{n} V(\vec{r}_{i})$ was the potential due to all the charges except \vec{r}_{i} . But while

writing the continuous form, we wrote, $W = \frac{1}{2} \iint P V d 2$

where Visis a continuous function, a ful potential. For continuous distribution, there is no problem, since the amount of charge right at \vec{p} is vanishingly small and the contribution to potential is zono. But for point charges, we should use the first equation.

Energy of uniformly charged spherical shell

$$W = \frac{1}{2} \int \sigma V dA = \frac{1}{2} \int \sigma \cdot \frac{1}{4\pi\kappa} \cdot \frac{9}{R} \cdot da$$

Swiface

 $= \frac{9}{8\pi\kappa} \cdot \frac{9}{R} \cdot \int \sigma' da = \frac{9^2}{8\pi\kappa} \cdot \frac{9}{R} \cdot da$

In field method,

$$W = \frac{40}{2} \iint E^2 d2 = \frac{40}{2} \left[\int E^2 d2 + \int E^2 d2 \right]$$
All opace inside ordere

$$= \int \frac{q^2}{16\pi^2 C^2} \cdot \frac{1}{84} r^2 \sin\theta \, dr \, d\theta \, dd$$

$$= \frac{q^2}{16\pi^2 C^2} \times 4\pi \int \frac{1}{r^2} \, dr$$

$$= \frac{q^2}{16\pi^2 C^2} \times 4\pi \int \frac{1}{r^2} \, dr$$

$$= \frac{9^2}{8116 R}$$

Energy for solid ophere of radius R and charge &

$$E_{in} = \frac{9}{36.0} \text{ Y} \qquad E_{out} = \frac{9R^{3}}{36.02}$$

$$V = \frac{6}{2} 41 \int \frac{92}{96.2} r^{2} dr + \frac{6}{2} 411 \int \frac{9^{2}p^{6}}{96.2} \frac{1}{r^{4}} r^{2} dr$$

$$= \frac{4116.9^{2}}{186.2} \int \frac{4116.9^{2}p^{6}}{186.2} \left[\frac{1}{R}\right]$$

$$= \frac{4116.9^{2}}{186.2} \left[\frac{R^{3}}{45} + R^{5}\right]$$

Superposition principle for energy

What = $\frac{E_2}{2} \int E^2 d2 = \frac{E_1}{2} \int (E_1 + E_2)^2 d2$ = $\frac{E_2}{2} \int (E_1^2 + E_2^2 + 2E_1 \cdot E_2) d2$ = $W_1 + W_2 + E_0 \int E_1 \cdot E_2 d2$

So, superposition principle doesn't hold.