LECTURE NO 19

Electrostatics

TOPIC COVERED

Energy density in the electrostatic field

To determine the energy present in an assembly of charges, we must first determine the amount of work necessary to assemble them. Suppose we wish to position three point charges Q_1 , Q_2 , and Q_3 in an initially empty space shown shaded in Figure 4.22. No work is required to transfer Q_1 from infinity to P_1 because the space is initially charge free and there is no electric field [from eq. (4.59), W = 0]. The work done in transferring Q_2 from infinity to P_2 is equal to the product of Q_2 and the potential V_{21} at P_2 due to Q_1 . Similarly, the work done in positioning Q_3 at P_3 is equal to $Q_3(V_{32} + V_{31})$, where V_{32} and V_{31} are the potentials at P_3 due to Q_2 and Q_1 , respectively. Hence the total work done in positioning the three charges is

$$W_E = W_1 + W_2 + W_3 = 0 + Q_2 V_{21} + Q_3 (V_{31} + V_{32})$$
(4.84)

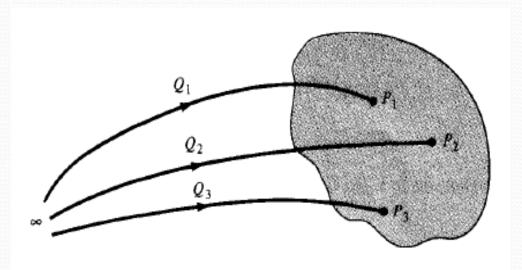
If the charges were positioned in reverse order,

$$W_E = W_3 + W_2 + W_1 = 0 + Q_2 V_{23} + Q_1 (V_{12} + V_{13})$$
(4.85)

where V_{23} is the potential at P_2 due to Q_3 , V_{12} and V_{13} are, respectively, the potentials at P_1 due to Q_2 and Q_3 . Adding eqs. (4.84) and (4.85) gives

$$2W_E = Q_1(V_{12} + V_{13}) + Q_2(V_{21} + V_{23}) + Q_3(V_{31} + V_{32})$$

= $Q_1V_1 + Q_2V_2 + Q_3V_3$



or

$$W_E = \frac{1}{2} (Q_1 V_1 + Q_2 V_2 + Q_3 V_3) \tag{4.86}$$

where V_1 , V_2 , and V_3 are total potentials at P_1 , P_2 , and P_3 , respectively. In general, if there are n point charges, eq. (4.86) becomes

$$W_E = \frac{1}{2} \sum_{k=1}^{n} Q_k V_k \qquad \text{(in joules)}$$
 (4.87)

If, instead of point charges, the region has a continuous charge distribution, the summation in eq. (4.87) becomes integration; that is,

$$W_E = \frac{1}{2} \int \rho_L V \, dl \qquad \text{(line charge)} \tag{4.88}$$

$$W_E = \frac{1}{2} \int \rho_S V \, dS \qquad \text{(surface charge)} \tag{4.89}$$

$$W_E = \frac{1}{2} \int \rho_{\nu} V \, d\nu \qquad \text{(volume charge)} \tag{4.90}$$

Since $\rho_{\nu} = \nabla \cdot \mathbf{D}$, eq. (4.90) can be further developed to yield

$$W_E = \frac{1}{2} \int_{v} (\nabla \cdot \mathbf{D}) \, V \, dv \tag{4.91}$$

But for any vector **A** and scalar *V*, the identity

$$\nabla \cdot V \mathbf{A} = \mathbf{A} \cdot \nabla V + V(\nabla \cdot \mathbf{A})$$

or

$$(\nabla \cdot \mathbf{A})V = \nabla \cdot V\mathbf{A} - \mathbf{A} \cdot \nabla V \tag{4.92}$$

holds. Applying the identity in eqs. (4.92) to (4.91), we get

$$W_E = \frac{1}{2} \int_{V} (\nabla \cdot V \mathbf{D}) \, dv - \frac{1}{2} \int_{V} (\mathbf{D} \cdot \nabla V) \, dv \tag{4.93}$$

By applying divergence theorem to the first term on the right-hand side of this equation, we have

$$W_E = \frac{1}{2} \oint_{S} (V\mathbf{D}) \cdot d\mathbf{S} - \frac{1}{2} \int_{V} (\mathbf{D} \cdot \nabla V) \, dV \tag{4.94}$$

From Section 4.9, we recall that V varies as 1/r and D as $1/r^2$ for point charges; V varies as $1/r^2$ and D as $1/r^3$ for dipoles; and so on. Hence, VD in the first term on the right-hand side of eq. (4.94) must vary at least as $1/r^3$ while dS varies as r^2 . Consequently, the first integral in eq. (4.94) must tend to zero as the surface S becomes large. Hence, eq. (4.94) reduces to

$$W_E = -\frac{1}{2} \int_{v} (\mathbf{D} \cdot \nabla V) \, dv = \frac{1}{2} \int_{v} (\mathbf{D} \cdot \mathbf{E}) \, dv \tag{4.95}$$

and since $\mathbf{E} = -\nabla V$ and $\mathbf{D} = \varepsilon_0 \mathbf{E}$

$$W_E = \frac{1}{2} \int \mathbf{D} \cdot \mathbf{E} \, dv = \frac{1}{2} \int \varepsilon_0 E^2 \, dv$$
 (4.96)

Assignment

Derive the expression for the energy density in the electrostatic field is given by