

Electrostatic Energy of Point Charges: Self-Energy and Interaction Energy [lam8]

We have noted in [ln5] that the expression,

$$U = \frac{1}{2} \frac{1}{4\pi\epsilon_0} \int d^3x \int d^3x' \frac{\rho(\mathbf{x})\rho(\mathbf{x}')}{|\mathbf{x} - \mathbf{x}'|} = \frac{1}{2} \int d^3x \rho(\mathbf{x})\Phi(\mathbf{x}), \quad (1)$$

for the potential energy of a charge distribution $\rho(\mathbf{x})$ is not equivalent to the expression,

$$U_{\text{int}} = \frac{1}{2} \frac{1}{4\pi\epsilon_0} \sum_{i \neq j} \frac{q_i q_j}{|\mathbf{x}_i - \mathbf{x}_j|}, \quad (2)$$

for the potential interaction energy of an array of discrete charges q_i . The former expression converted into an integral over energy density $w(\mathbf{x})$ as demonstrated in [ln5],

$$U = \frac{1}{2} \int d^3x \rho(\mathbf{x})\Phi(\mathbf{x}) = \frac{1}{2} \epsilon_0 \int d^3x |\mathbf{E}(\mathbf{x})|^2 = \int d^3x w(\mathbf{x}), \quad (3)$$

is manifestly non-negative, whereas the latter expression can be negative (e.g. for two point charges of opposite sign).

We have also shown in [ln5] how the the integral expressions in (3) can be split into the sum of self-energy and interaction energy.

The energy density generated by the electric field of two point charges is

$$\begin{aligned} w(\mathbf{x}) &= \frac{1}{2} \epsilon_0 \left[\frac{q_1}{4\pi\epsilon_0} \frac{\mathbf{x} - \mathbf{x}_1}{|\mathbf{x} - \mathbf{x}_1|^3} + \frac{q_2}{4\pi\epsilon_0} \frac{\mathbf{x} - \mathbf{x}_2}{|\mathbf{x} - \mathbf{x}_2|^3} \right]^2 \\ &= \frac{1}{32\pi^2\epsilon_0} \left[\frac{q_1^2}{|\mathbf{x} - \mathbf{x}_1|^4} + \frac{q_2^2}{|\mathbf{x} - \mathbf{x}_2|^4} + 2q_1q_2 \frac{(\mathbf{x} - \mathbf{x}_1) \cdot (\mathbf{x} - \mathbf{x}_2)}{|\mathbf{x} - \mathbf{x}_1|^3 |\mathbf{x} - \mathbf{x}_2|^3} \right]. \quad (4) \end{aligned}$$

The first two terms represent, upon integration, the electrostatic self-energies, which are divergent and not included in U_{int} as constructed in (2).

Divergent charge densities or energy densities do not, in general, produce infinite self-energies. The problem here is caused by having finite charges concentrated in points, which yields strong (non-integrable) divergences.

The last term in (4) represents U_{int} and can be transformed (upon integration) as follows [lex120]:

$$U_{\text{int}} = \frac{q_1 q_2}{16\pi^2\epsilon_0} \int d^3x \frac{(\mathbf{x} - \mathbf{x}_1) \cdot (\mathbf{x} - \mathbf{x}_2)}{|\mathbf{x} - \mathbf{x}_1|^3 |\mathbf{x} - \mathbf{x}_2|^3} = \frac{1}{4\pi\epsilon_0} \frac{q_1 q_2}{|\mathbf{x}_1 - \mathbf{x}_2|}.$$