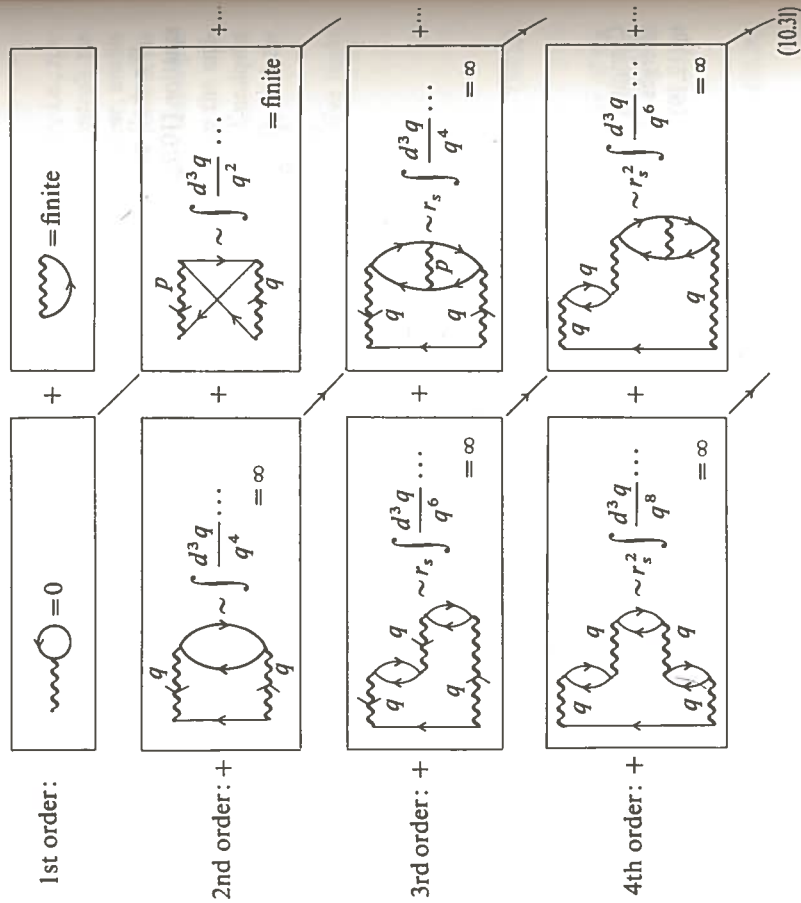


A look at the higher orders shows the same behaviour: most of the diagrams diverge. This appears to be a slightly discouraging start. However, the situation is saved by the partial summation method as follows: First of all, we observe that each term in the self-energy is proportional to some power of  $r_s$  ( $\sim 1/k_F$ —see §10.3). To see this, consider the second-order self-energy term in (9.66). To make the integral dimensionless, express quantities in units of  $k_F$ , i.e., set  $q = q'/k_F$ ,  $\epsilon_l = \epsilon_l'/k_F^2$ , etc. Then we find that (9.66)  $\sim k_F^3 \times k_F^3 \times k_F^4 \times k_F^{-2} = 1$ , so that (9.66) has no explicit dependence on  $k_F$ . Now look at a third-order term—e.g., the fifth or sixth diagram on the right of (10.8). After doing the frequency integrals, we find, compared with the second-order term, an extra  $V_q$  ( $\sim k_F^2$ ), an extra energy denominator ( $\sim k_F^{-2}$ ) and an extra integral over  $k$  ( $\sim k_F^3$ ), so that we have an extra factor  $k_F^{-1} \sim r_s$ . Thus all third-order terms are  $\sim r_s$ . In general, any  $n$ th-order term  $\sim r_s^{n-2}$ .

Next, we arrange the diagrams according to degree of divergence (= number of factors  $q^2$  in the denominator of the integrand) and dependence on  $r_s$ :



The first-order bubble has  $q=0$ , so it vanishes by (10.28); that is, the bubble is cancelled by the positive background. We see that the digrams can be arranged so that in each order of divergence they form a power series in  $r_s$  (see along diagonal lines). Thus, for small  $r_s$ , i.e., high-density limit, the dominant infinite terms are just those of lowest order in  $r_s$ . Hence in the high density limit, the self-energy series is just the sum

(10.32)

i.e., the sum over all diagrams of the repeated pair-bubble or 'ring' type. The remarkable thing is that this sum over an infinite number of infinite terms can actually be carried out, and it gives a finite result! Approximation (10.32) for the self-energy is called the 'random phase approximation' or 'RPA', for historical reasons.

The sum over rings is easy. Factoring out a free propagator from each diagram in (10.32) (this is not quite straightforward for the oyster, since it is a special case, by the rules of Table 9.1; however, it is legitimate) gives

(10.33)

The double wiggle is the 'effective interaction' in RPA, which was mentioned in §4.9, and interpreted as a 'screened' interaction between two particles:

(1) (2) (3)

(10.34)