Every distinct state of a binary alloy system on N sites is contained in the symbolic product of N factors:

$$(A_1 + B_1)(A_2 + B_2)(A_3 + B_3) \cdots (A_N + B_N)$$
, (19)

in analogy to (4). The average composition of a binary alloy is specified conventionally by the chemical formula  $A_{1-x}B_x$ , which means that out of a total of N atoms, the number of A atoms is  $N_A = (1-x)N$  and the number of B atoms is  $N_B = xN$ . Here x lies between 0 and 1.

The symbolic expression

$$(A + B)^{N} = \sum_{t=0}^{N} \frac{N!}{(N-t)! \ t!} A^{N-t} B^{t}$$
 (20)

is analogous to the result (12). The coefficient of the term in  $A^{N-t}$  B' gives the number g(N,t) of possible arrangements or states of N-t atoms A and t atoms B on N sites:

$$g(N,t) = \frac{N!}{(N-t)! \, t!} = \frac{N!}{N_A! \, N_B!} \,, \tag{21}$$

which is identical to the result (15) for the spin model system, except for notation.

## Sharpness of the Multiplicity Function

We know from common experience that systems held at constant temperature usually have well-defined properties; this stability of physical properties is a major prediction of thermal physics. The stability follows as a consequence of the exceedingly sharp peak in the multiplicity function and of the steep variation of that function away from the peak. We can show explicitly that for a very large system, the function g(N,s) defined by (15) is peaked very sharply about the value s = 0. We look for an approximation that allows us to examine the form of g(N,s) versus s when  $N \gg 1$  and  $|s| \ll N$ . We cannot look up these values in tables: common tables of factorials do not go above N = 100, and we may be interested in  $N \approx 10^{20}$ , of the order of the number of atoms in a solid specimen big enough to be seen and felt. An approximation is clearly needed, and a good one is available.

It is convenient to work with  $\log g$ . Except where otherwise specified, all logarithms are understood to be  $\log \operatorname{base} e$ , written here as  $\log$ . The international standard usage is  $\ln \operatorname{for} \log \operatorname{base} e$ , but it is clearer to write  $\log \operatorname{when}$  there is no ambiguity whatever. When you confront a very, very large number such as

