

**Figure 3-4**

In this figure, we show the curves relevant to BALs we can observe, *e.g.* C IV  $\lambda 1549$ , Si IV  $\lambda 1397$ , O VI  $\lambda 1034$ , etc. The  $1s^2 2s^1$  isoelectronic sequence (see chapter 9) is shown in solid lines, the  $1s^2 2s^2 2p^6 3s^1$  sequence is shown as dotted lines, and the curve for neutral hydrogen is shown as a dashed line. Since we might see Fe III  $\lambda 1122$  in QSOs with the lower ionization BALs, Fe<sup>+3</sup> is also shown as a dash-dot line. CLOUDY has no parameters for Phosphorus, so we have made a rough estimate for P<sup>+4</sup> by averaging the curves for Si<sup>+4</sup> and S<sup>+4</sup> which have similar values of  $U_p$  (see figure 3-3). The peaks of these ions, in order of increasing Z, are:  $U_p(\text{C}^{+3}) = -1.86$ ,  $U_p(\text{N}^{+4}) = -1.29$ ,  $U_p(\text{O}^{+5}) = -0.64$ ,  $U_p(\text{Al}^{+2}) = -2.68$ ,  $U_p(\text{Si}^{+3}) = -2.08$ ,  $U_p(\text{S}^{+5}) = -1.27$ , and the rough estimate for the peak for Phosphorus is  $U_p(\text{P}^{+4}) = -1.59$ .

$N(\text{H}) = 10^{18}$  was used, and both sides of the slab were averaged.