

Erratum: Energy relaxation and the quasiequation of state of a dense two-temperature nonequilibrium plasma
[Phys. Rev. E 58, 3705 (1998)]

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We wish to correct a numerical mistake in Table II where the numbers have to be rescaled by a multiplicative factor. Here we present revised values, where we have also used a more recent local-field correction, viz., see Phys. Rev. B **62**, 16536 (2000), in the electron response functions. The ion-ion static structure factor for Al at the melting point has also been recalculated using an Ashcroft-type pseudopotential described in Phys. Rev. B **28**, 1701 (1983). For higher electron temperatures relevant to the nonequilibrium problem, the pseudopotentials given in Table I were used as before. The revised values of the coupling constants calculated using the Fermi golden rule (FGR) and the Spitzer model (S) are within about a factor of 3. The coupled-mode (CM) results are an order of magnitude smaller. Hence the main conclusions of the paper remain unchanged. The corrected table is given below [1].

TABLE I. Calculated values of the electron-ion coupling constant \bar{g}_{cc} for Al at normal density and with $T_i=0.08126$ eV. Results from the Spitzer (S), Fermi golden rule (FGR), and coupled mode (CM) calculation are given in Watts/Kelvin/cubic meter.

| T (eV) | S/ 10^{19} | FGR/ 10^{18} | CM/ 10^{17} |
|----------|--------------|----------------|---------------|
| 3.0 | 0.1998 | 0.6100 | 0.5469 |
| 4.0 | 0.2216 | 0.6386 | 0.6064 |
| 5.0 | 0.2360 | 0.6634 | 0.6375 |
| 7.0 | 0.2504 | 0.6897 | 0.6358 |
| 9.0 | 0.2390 | 0.7805 | 0.6132 |
| 10.0 | 0.2211 | 0.8679 | 0.6029 |
| 15.0 | 0.2079 | 1.0206 | 0.7942 |
| 20.0 | 0.2071 | 1.2962 | 0.9150 |
| 25.0 | 0.2267 | 1.1753 | 1.0697 |
| 30.0 | 0.2493 | 1.4276 | 1.3099 |
| 35.0 | 0.2829 | 1.8898 | 1.5673 |
| 40.0 | 0.3191 | 2.0012 | 1.3346 |

[1] More details are given at the website: <http://nrcphy1.phy.nrc.ca/ims/qp/chandre/Lst/>.