ERRATUM

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GALLISTEL, C. R. AND A. J. DAVIS. Affinity for the dopamine D_2 receptor predicts neuroleptic potency in blocking the reinforcing effect of MFB stimulation. PHARMACOL BIOCHEM BEHAV 19(5) 867–872, 1983.

The authors wish to correct the following errors. The references cited in the legends for Figs. 1 and 2 were incorrectly numbered. The complete legends, with the correct reference numbers, are reproduced below.

FIG. 1. The logarithm of the neuroleptic dose required to block sustained self-stimulation plotted against the logarithm of affinity for the dopamine D₂ receptor, as measured in two in vitro studies. The number beside a point gives the number of rats for which that was the required dose. The best fitting linear regressions (solid lines) were computed from the logarithmic mean of the required doses of a given drug. Both correlations are significant beyond the .005 level. Abbreviations: Ben=benperidol; Chlor=chlorpromazine; Cloz= clozapine; Halo=haloperidol; Meto=metoclopromide; Pim= pimozide; Prom=promazine; Spiro=spiroperidol; Thio=thioridizine. A. Affinity measured by Creese, Burt, and Snyder [2] using ³H-halopridol as the radioactive ligand and calf striatal tissue. (K₁ is proportional to IC₅₀, which is the concentration of a drug required to displace 50% of the stereospecifically bound ligand.) B. Affinity measured by Leysen, Gommeren, and Laduron [16] using 3Hspiroperidol and rat striatal tissue. (pIC₅₀ = $-\log IC_{50}$. *In calculating the linear regression, the sign of the numbers on the abscissa was changed where necessary to make them increase in value from left to right.)

FIG. 2. The mean logarithm of the neuroleptic dose required to block sustained self-stimulation plotted against the logarithm of the drugs' affinities for various aminergic receptors. (In each case the high affinity is placed at the left end of the abscissa. In calculating the linear regression, the sign of the numbers on the abscissa was changed where necessary to make them increase in value from left to right. None of the correlations approaches statistical significance. For definitions of IC₅₀, pIC₅₀, and K_i see caption to FIG. 1. Abbreviations: B=benperidol; Ch=chlorpromazine; Cl=clozapine; H=haloperidol; M=metoclopromide; Pi=pimozide; Pr=promazine; S=spiroperidol; Th=thioridizine.) A. Affinity data for the D₁ dopamine receptor from [13] with the ³H-cis(2)-flupenthixol ligand and rat striatal tissue. B. Affinity data for the D₃ dopamine receptor from [2] with the ³H-dopamine ligand and calf striatal tissue. C. Affinity data on the α -adrenergic receptor from [18] with the ³H-WB-4101 ligand and rat whole brain. D. Affinity data for the S₂ serotonin receptor from [15] with the 3H-spiroperidol ligand and rat frontal cortex. E. Affinity data for the S_1 serotonin receptor in the cortex from [17] with the ³H-serotonin ligand and rat cortical tissue. F. Affinity data for the S₁ serotonin receptor in the hippocampus from [15] with the ³Hserotonin ligand and rat hippocampal tissue.