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Erratum

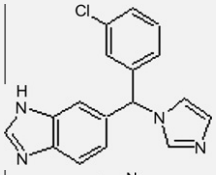
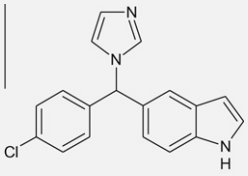
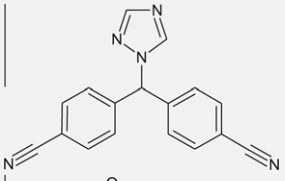
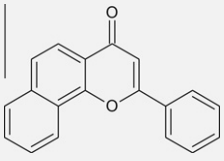
Erratum to "Pharmacophore modeling strategies for the development of novel nonsteroidal inhibitors of human aromatase (CYP19)" [Bioorg. Med. Chem. Lett. 20 (2010) 3050]

Yagmur Muftuoglu [a], Gabriela Mustata [b,*]

[a] Departments of Biophysics and Chemical Biomolecular Engineering, Johns Hopkins University, Baltmore, MD 21218, United States
[b] Department of Computational Biology, School of Medicine, University of Pittsburgh, Pittsburgh, PA 15213, United States

The publisher regrets that Table 2 appeared incorrectly. The correct table appears below.

Table 2
Active test set of 36 active and structurally diverse aromatase inhibitors

#	Compound name	Chemical structure	IC_{50}
1	Liarozole[57]		5 nM
2	5-[(4-Chlorophen-yl) (1*H*-imidazol-1-yl)methyl]-1*H*-indole[58]		9 nM
3	Letrozole[14]		11.5 nM
4	7,8-Benzoflavone[61]		70 nM

(*continued on next page*)

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* Corresponding author. Tel.: +1 412 648 7799; fax: +1 412 648 3163.
E-mail address: gmustata@pitt.edu (G. Mustata).

doi:10.1016/j.bmcl.2010.06.115

Table 2 (*continued*)

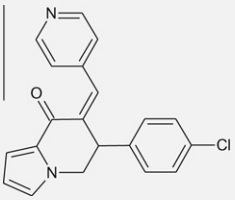
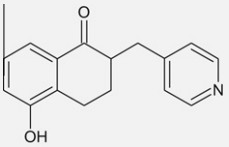
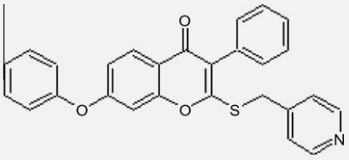
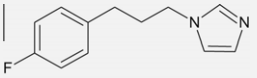
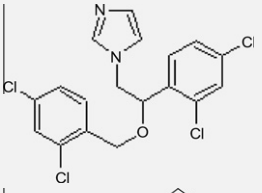
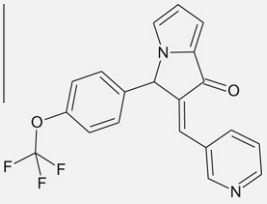
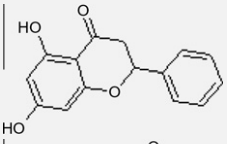
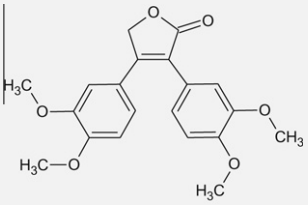
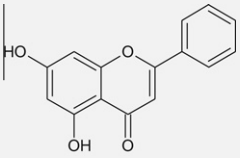
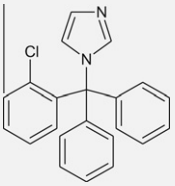
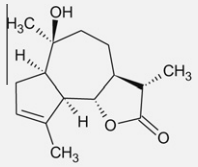
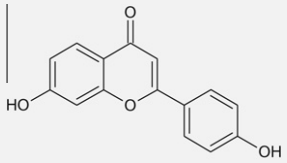
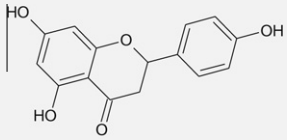
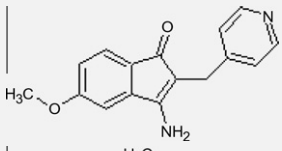
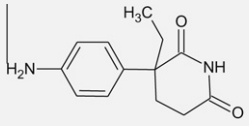
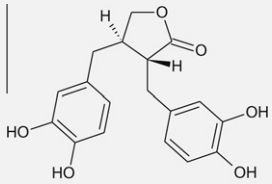
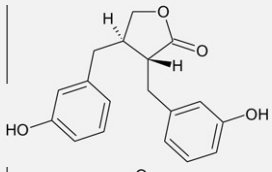
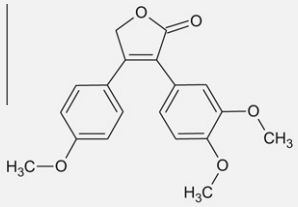
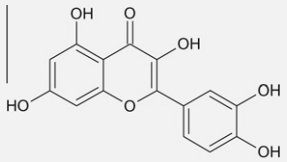
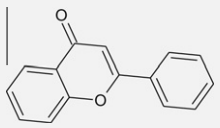
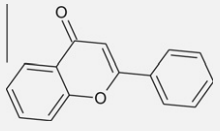
#	Compound name	Chemical structure	IC_{50}
5	MR 20492[63]		200 nM
6	35075-39-7[65]		195 nM
7	5j[45]		210 nM
8	4F[51]		310 nM
9	Miconazole[60]		600 nM
10	MR 16089[48]		650 nM
11	Apigenin[59]		900 nM
12	TM-8[54]		1.0 μM
13	Chrysin[54,59]		1 μM
14	Clotrimazole[60]		1.8 μM

Table 2 (*continued*)

#	Compound name	Chemical structure	IC_{50}
15	11β,13-Dihydro-10-*epi*-8-deoxycumam-brin B[62]		2.0 μM
16	7,4′-Dihydroxyflavone[64]		2.0 μM
17	Naringenin[54]		2.9 μM
18	MR 20814[66]		3.5 μM
19	Aminoglutethimide[67]		5.2 μM
20	4,4′-Dihydroxy enterolactone[68]		6 μM
21	Enterolactone[68]		6 μM
22	TM-7[54]		6.9 μM
23	Quercetin[50]		12 μM
24	Flavone[61]		8.0 μM
25	Flavanone[59]		8.7 μM

(*continued on next page*)

Table 2 (*continued*)

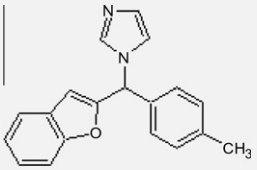
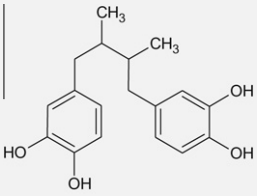
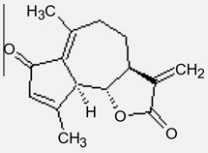
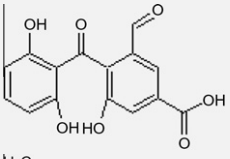
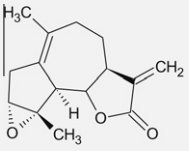
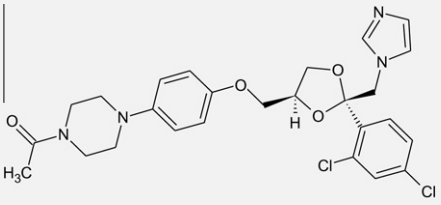
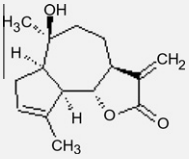
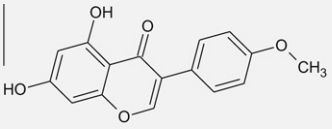
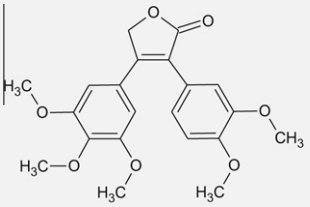
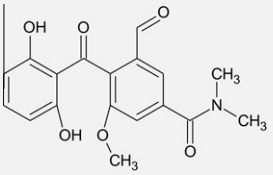
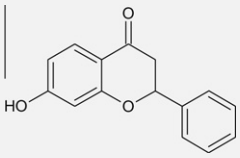
#	Compound name	Chemical structure	IC_{50}
26	4C[69]		8.8 μM
27	Nordihydroguaiaretic Acid[68]		11 μM
28	Dehydroleucodin[62]		15 μM
29	TAN-931[70]		17.2 μM
30	Ludartin[62]		55 μM
31	Ketoconazole[60]		60 μM
32	10-Epi-8-Deoxy-cumambrin B[62]		7 μM
33	Biochanin A[59]		10.2 μM
34	TM-9[54]		13.7 μM
35	2s[70]		17.2 μM

Table 2 (*continued*)

#	Compound name	Chemical structure	IC_{50}
36	7-Hydroxyflavone[59]		30.5 μM