

## **Drug Metabolism and Disposition**

### **A PREDICTIVE LIGAND-BASED BAYESIAN MODEL FOR HUMAN DRUG INDUCED LIVER INJURY**

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**Supplemental Table 4.** Mean physicochemical properties for Recently approved drugs from Prous database

Descriptor	Recently approved drugs from Prous database (N = 77)
ALogP	2.09 ± 3.49
Apol	16315.18 ± 9937.28
LogD	1.42 ± 3.52
MW	427.05 ± 280.31
Number of rotatable bonds	7.05 ± 7.56
Number of rings	3.44 ± 1.70
Number of aromatic rings	2.02 ± 1.21
Number of H bond acceptors	6.01 ± 6.73

Number of H bond donors	$2.37 \pm 3.28$
Molecular surface area	$413.89 \pm 264.25$
Molecular polar surface area	$110.85 \pm 133.18$
Wiener Index	$5843.43 \pm 17813.73$
Zagreb Index	$158.23 \pm 97.50$