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Table S1. Drug loss predicted by using 1D model for various drugs

Compound Name	MW (Da)	LogK*	D _p ** (m2/s)	Loss in PDMS at various flow rates ^{***} (%)			$\frac{C_{i,sl}^{*}}{C_{i,sl}}$ for various separation distance ^{\$\$}		
				1000 (ml/day)	10 (ml/day)	0.1 (ml/day)	0.1 mm	1 mm	10 mm
Cancer drugs									
Sunitinib	398	5.2	1.4E-14	5-25%	>50%	>50%	4.2	0.0	0.0
Doxorubicin	543	1.3	6.2E-15	<5%	<5%	5-25%	0.2	0.0	0.0
Vincristine	824	3.4	2.1E-15	<5%	5-25%	>50%	0.0	0.0	0.0
Cisplatin	298	-2.2	3.0E-14	<5%	<5%	<5%	16.2	0.0	0.0
Fluorouracil	130	-0.9	2.6E-13	<5%	<5%	<5%	63.5	0.0	0.0
Cardiovascular drugs									
epinephrine	183	-0.8	1.1E-13	<5%	<5%	<5%	45.9	0.0	0.0
Flecainide	414	3.8	1.3E-14	<5%	>50%	>50%	3.2	0.0	0.0
Isoproterenol	211	0.3	7.3E-14	<5%	<5%	5-25%	37.2	0.0	0.0
Metoprolol	267	1.8	3.9E-14	<5%	<5%	>50%	22.6	0.0	0.0
Verapamil	454	5.2	9.9E-15	5-25%	>50%	>50%	1.6	0.0	0.0

*Found

from (ref ¹)

**The Dp (diffusion coefficient of drugs in PDMS) was found by using following scaling relation as used previously ². We used paclitaxel diffusion coefficient found in this study for the scaling.

$$\mathsf{D}_{\mathsf{p},\,\mathsf{u}} = \mathsf{D}_{\mathsf{p},\,\mathsf{ptaxel}} \, (\overline{^{MW}_{Ptaxel}})^{-2.6}$$

Where, $D_{p, ptaxel}$ and MW_{ptaxel} are the diffusion coefficient and molecular weight of paclitaxel; $D_{p, u}$ and MW_{ptaxel} is the diffusion coefficient and molecular weight of the drug compound under analysis

***The loss was calculated for various flow rates mentioned for a microfluidic channel of 100µm x 200µm x 100mm (h x w x l) at t = 24 hr using equation 7.

$\frac{C_{i,sl}}{C_{i,sl}}$ was calculated using equation 11 for various separation distances (x) mentioned in the table.

References

- 1. D. S. Wishart, C. Knox, A. C. Guo, S. Shrivastava, M. Hassanali, P. Stothard, Z. Chang and J. Woolsey, *Nucleic Acids Res*, 2006, **34**, D668-672.
- 2. O. J. Karlsson, J. M. Stubbs, L. E. Karlsson and D. C. Sundberg, *Polymer*, 2001, **42**, 4915-4923.