Table S1. Structures and LSD1 inhibitory activities of compounds 1-35 with embedded structures.

	\mathbb{R}^1	\mathbb{R}^2	LSD1 IC ₅₀
			(µM)
1	4-OBn	Ņ	0.0098
		Z _Z N	
		Ö	
2	4-(6-F-pyridin-3-yl)	same as above	0.077
3	4-Br	same as above	0.035
4	4-(6-C1-pyridin-3-yl)	same as above	0.15
5	2-Br	same as above	>100
6	3-Br	same as above	>100
7	4-OBn	-Н	0.91
8	6-F-pyridin-3-yl	-Н	5.3
9	4-Br	-Н	25
10	4-Br	NH	0.064
		3 ₂ , N	
11	4-Br	3 ₂ N	0.62
		0	
12	4-Br	32, N	3.6
13	4-H	same as 10	0.19
14	4-H	same as 11	1.3
15	4-Br	NH	0.26
		ZZ-	
16	4-H	same as above	0.062
17	4-Br	-Bn	5.0
	l .		1

18	4-Br	-(4-OH)-Bn	94
19	4-Br	-(4-NH ₂)-Bn	>100
20	4-Br	-(4-NMe ₂)-Bn	71
21	4-Br	(pyridin-4-yl)-methyl	>100
22	4-Br	-COCH ₃	15.4
23	4-Br	-CONHPh	7.4
24	4-Br	=CH-Ph	15.6
25	4-Br	=CH-(4-Br-Ph)	1.2
26	4-Br	=CH-(2-OMe-Ph)	0.74
27	4-Br	=CH-(3-OMe-Ph)	0.90
28	4-Br	=CH-(2,3,4-OMe-Ph)	94
29	4-Br	=CH-(4-NMe ₂ -Ph)	11
30	4-Br	=CH-(pyridine-4-yl)	14
31	4-H	-NH ₂	5.8
32	4-H	yz N O NH	0.67
33	4-H	H N NH	84
34	4-Br	H NH	1.7
35	Ph	N N	36.4