Erratum

Spectrometrically monitored selection experiments: quantitative laser desorption mass spectrometry of small chemical libraries

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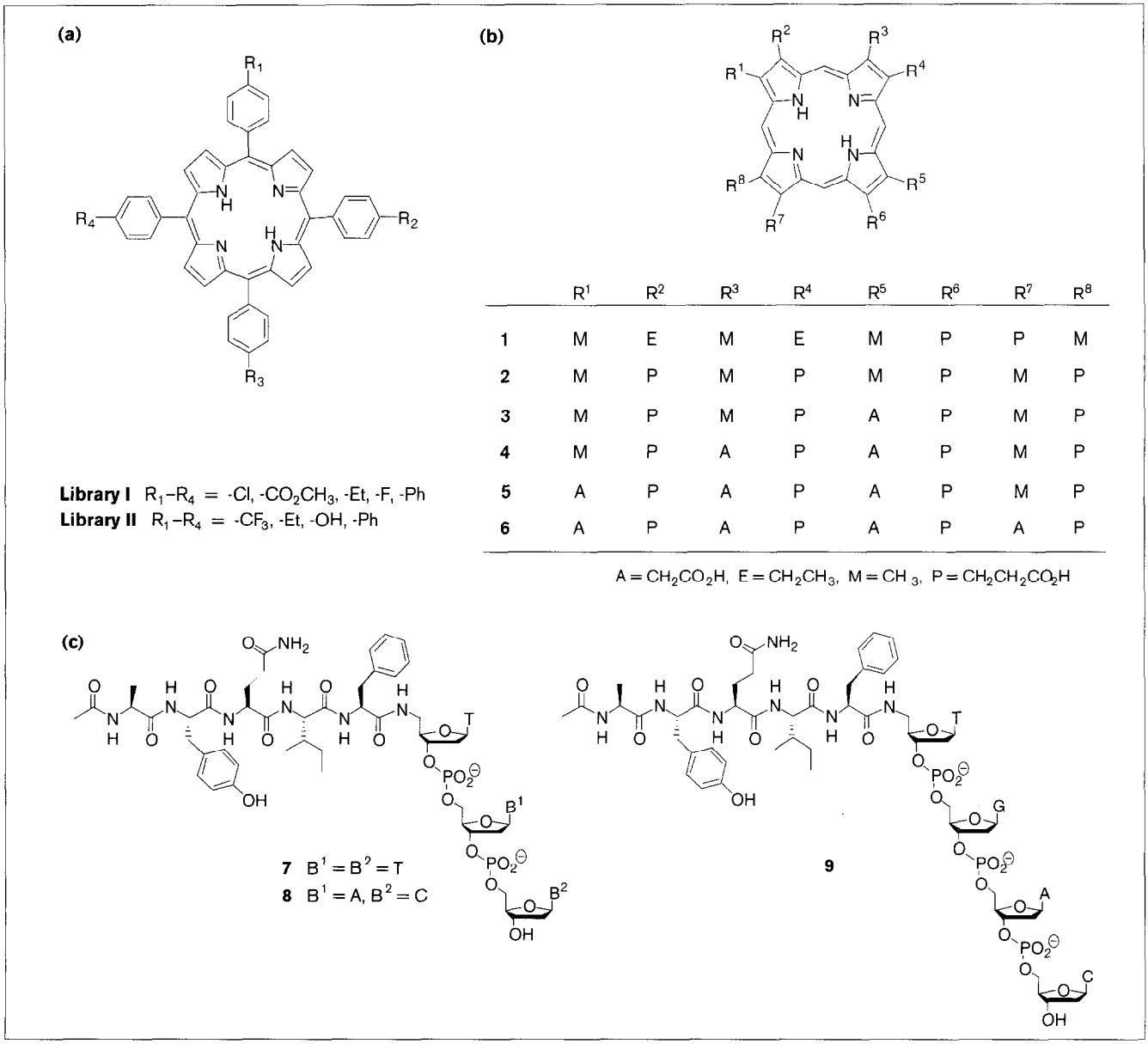
Chemistry & Biology January 1997, **4**:63–77

Errors were inadvertantly introduced into Figures 2 and 3 of this paper.
The figures are correctly reprinted below.

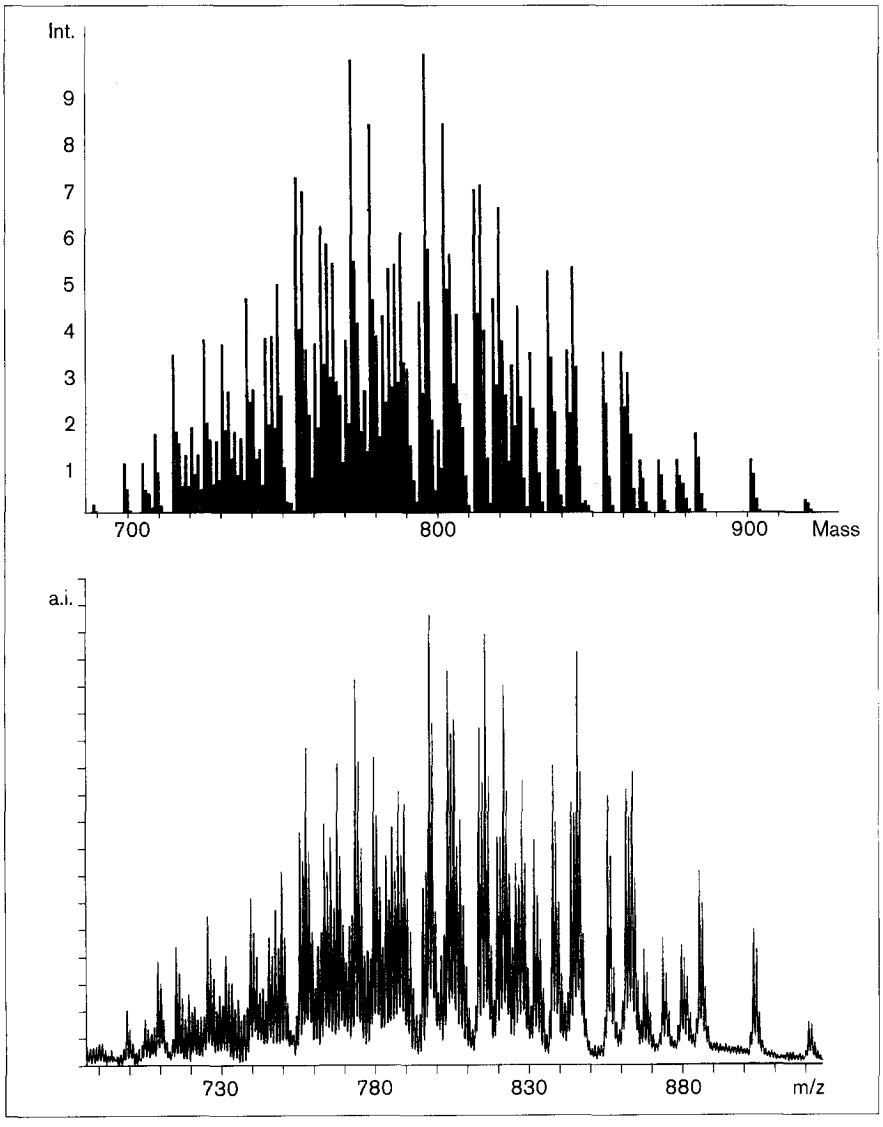
Figure 2



	R1	R2	R3	R4	R5	R6	R7	R8
1	M	E	M	E	M	P	P	M
2	M	P	M	P	M	P	M	P
3	M	P	M	P	A	P	M	P
4	M	P	A	P	A	P	M	P
5	A	P	A	P	A	P	M	P
6	A	P	A	P	A	P	A	P

Structural formulae of **(a)** tetraphenylporphyrin libraries, **(b)** porphyrin acids and **(c)** peptide–DNA hybrids employed in this work.

Figure 3



Mass spectrum of **Library I**. Top spectrum, predicted isotopically-resolved mass spectrum of **Library I** generated with the computer program MASP and bottom spectrum, experimental MALDI-TOF mass spectrum of **Library I**. This library contains 70 non-isobaric compounds, each of which is producing a set of isotope peaks. Peak intensity deviations from the predicted spectrum (top) are due to different reactivities of building blocks used in the combinatorial synthesis.