Electronic Supplementary Material (ESI) for Organic & Biomolecular Chemistry.

Supporting Information

Convergent synthesis of hydrophilic monomethyl dolastatin 10 based drug linkers for antibody drug conjugation

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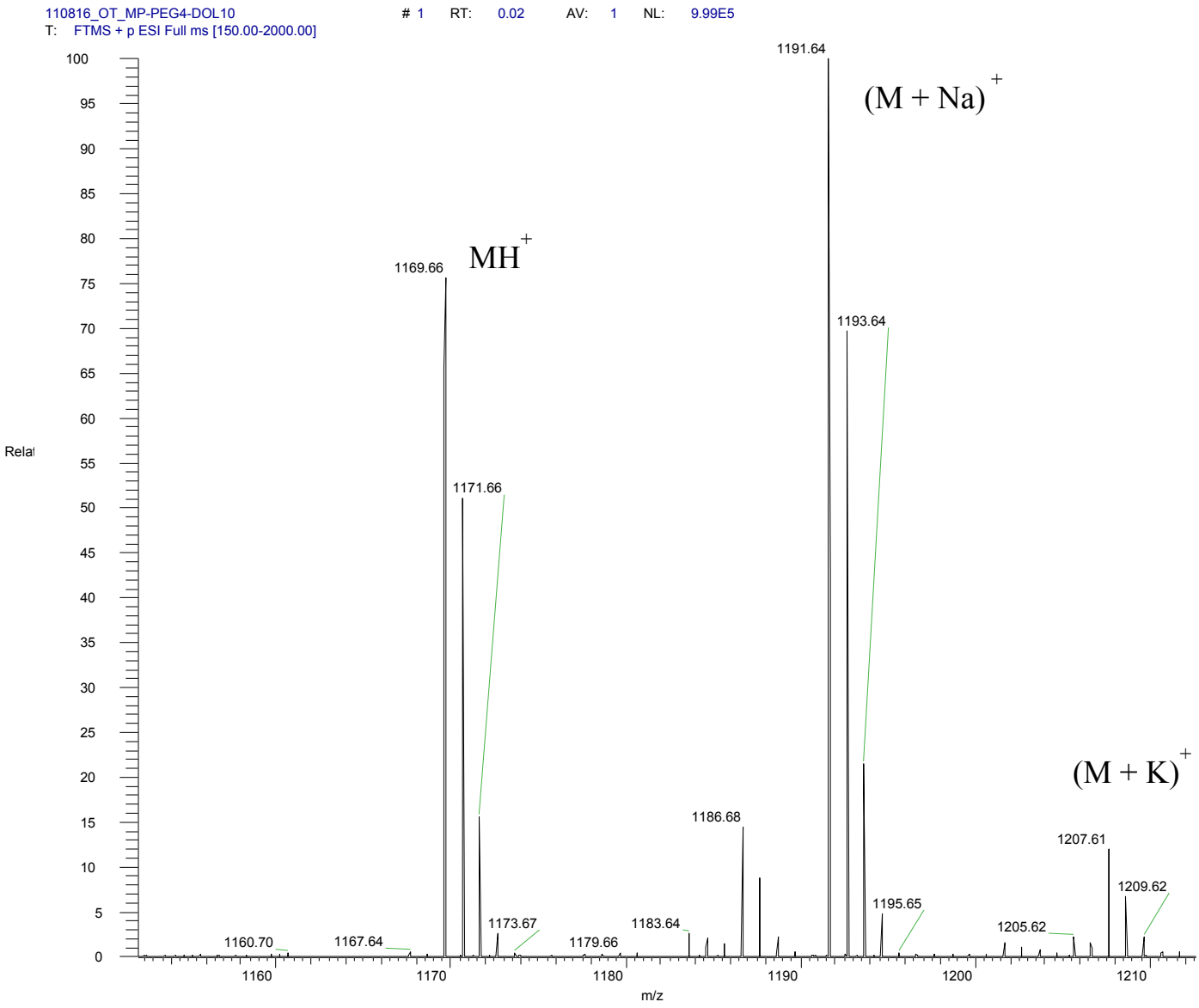


Figure S1. ESI mass spectrum of maleimido propyl $(PEG)_4$ dolastatin 10 drug linker with theoretical m/z of MH+ = 1169.65 (please note that in addition to protonated molecular ion peak, both sodium and potassium ion complexed molecular ions peaks were observed.).

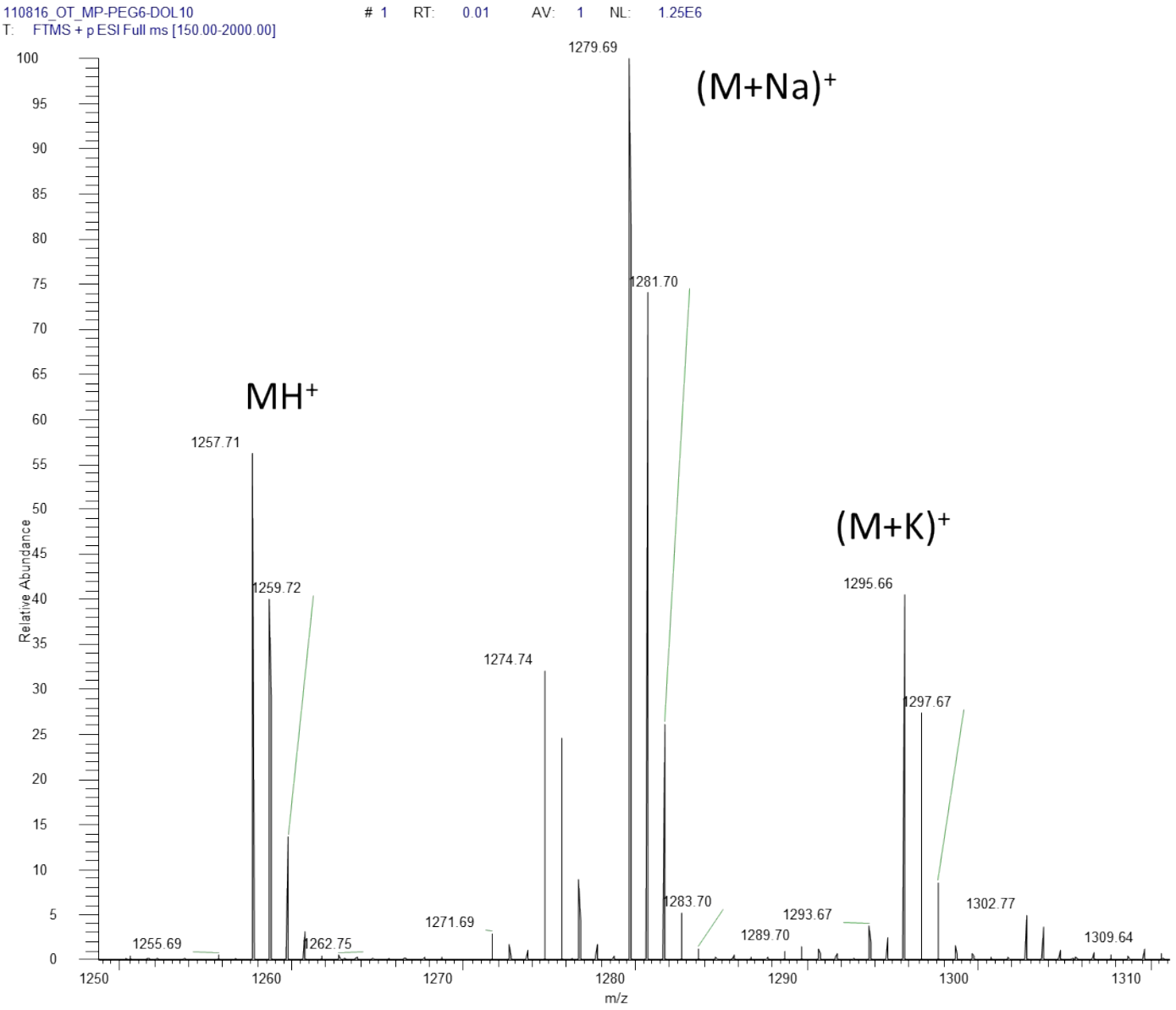


Figure S2. ESI mass spectrum of maleimido propyl $(PEG)_6$ dolastatin 10 drug linker with theoretical m/z of MH+ = 1257.70 (please note that in addition to protonated molecular ion peak, both sodium and potassium ion complexed molecular ions peaks were observed.).

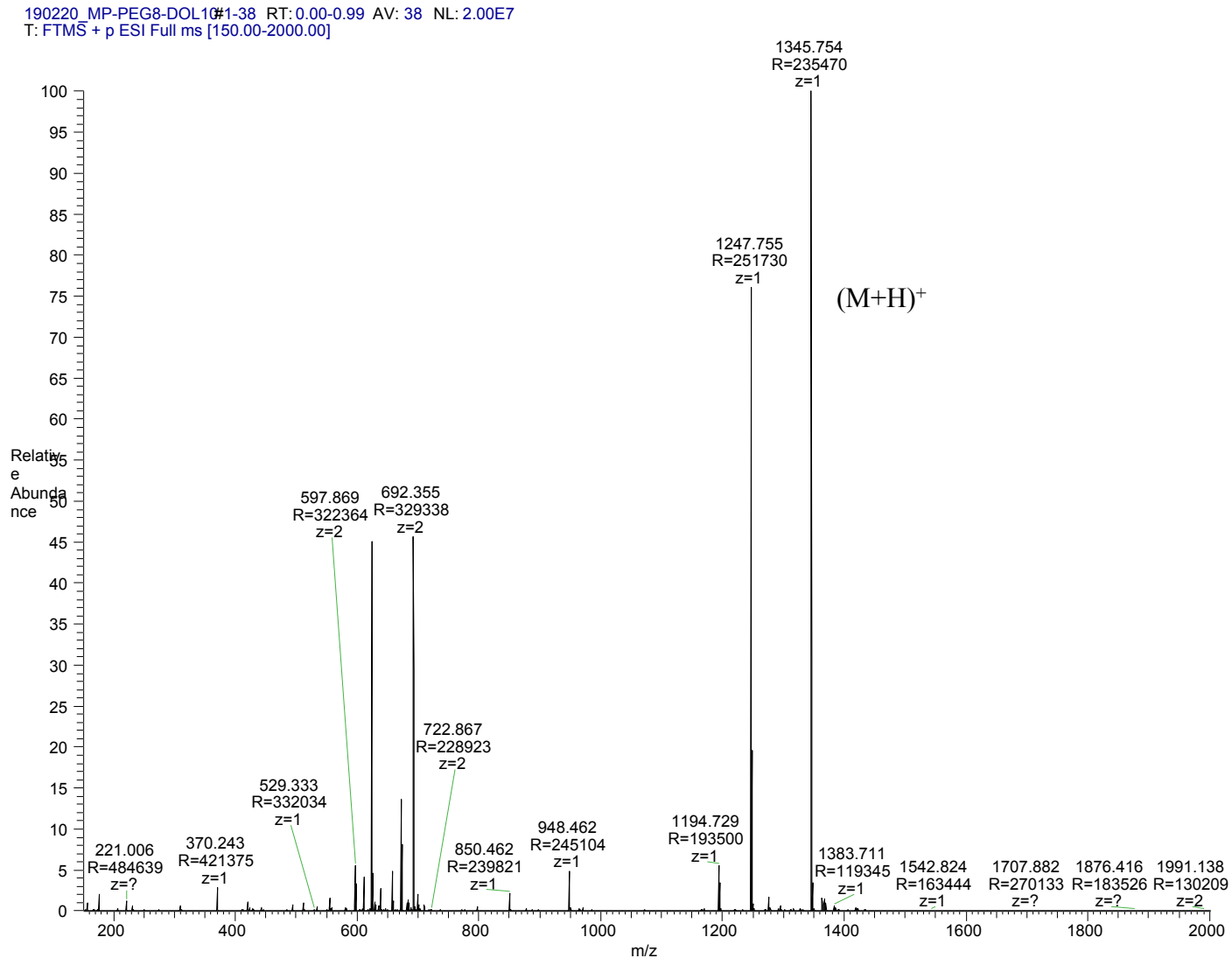


Figure S3. ESI mass spectrum of maleimido propyl $(PEG)_8$ dolastatin 10 drug linker (theoretical m/z of MH+ =1345.75 Dalton)

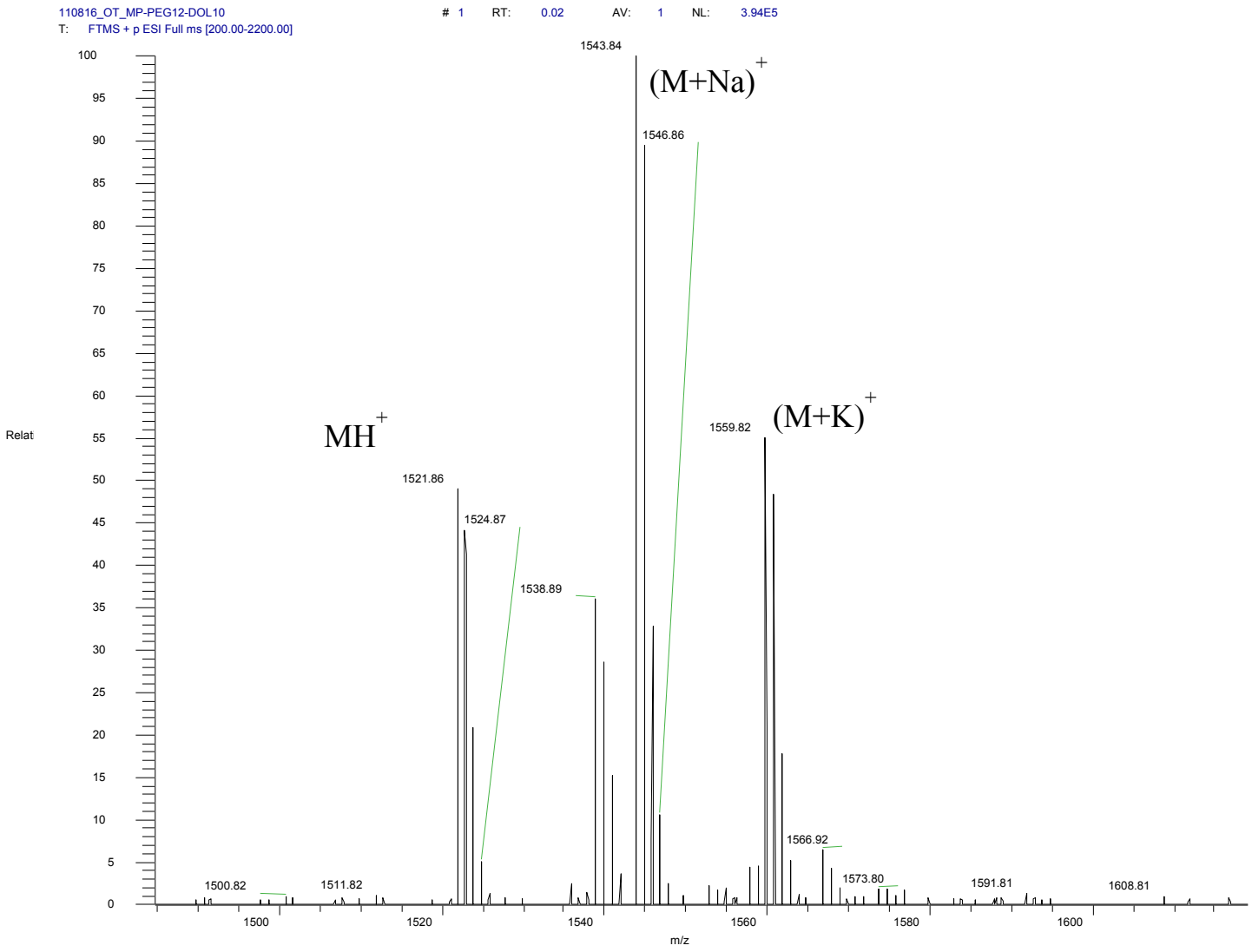


Figure S4. ESI mass spectrum of maleimido propyl $(PEG)_{12}$ dolastatin 10 drug linker with theoretical m/z of MH+ = 1521.86 (please note that in addition to protonated molecular ion peak, both sodium and potassium ion complexed molecular ions peaks were observed.).

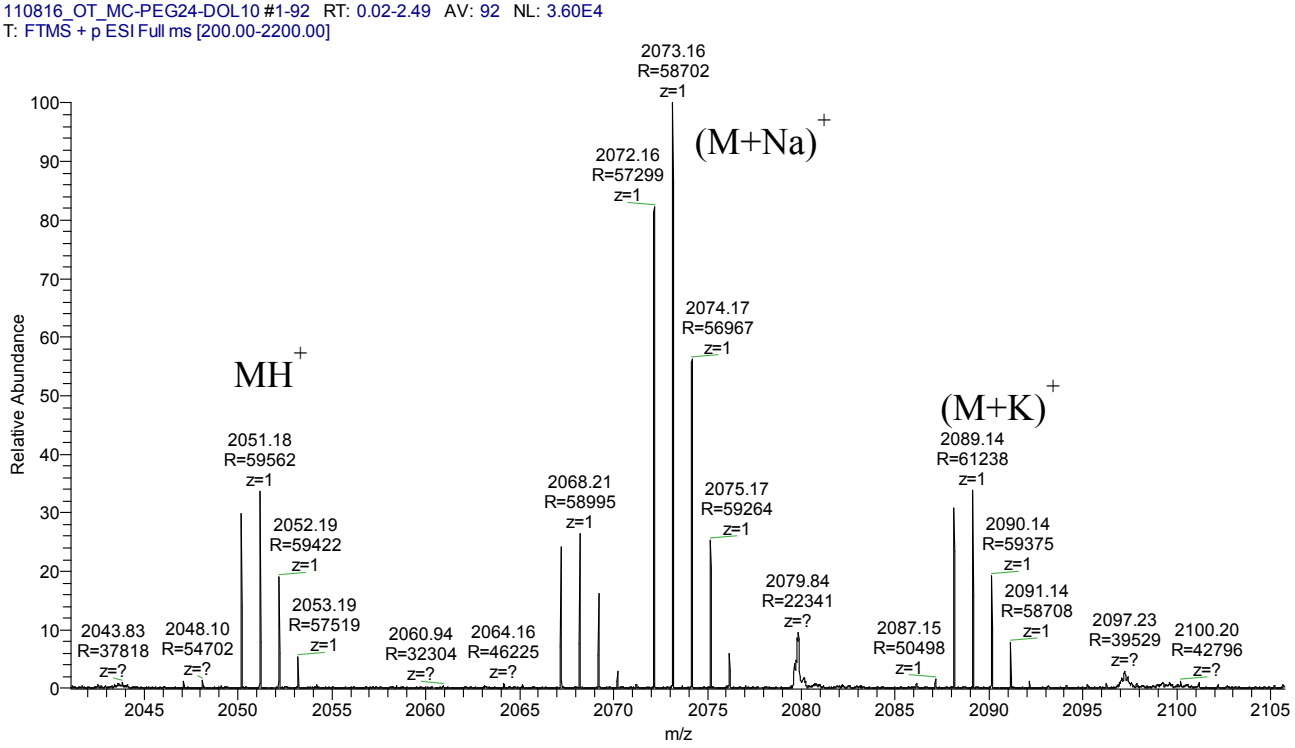
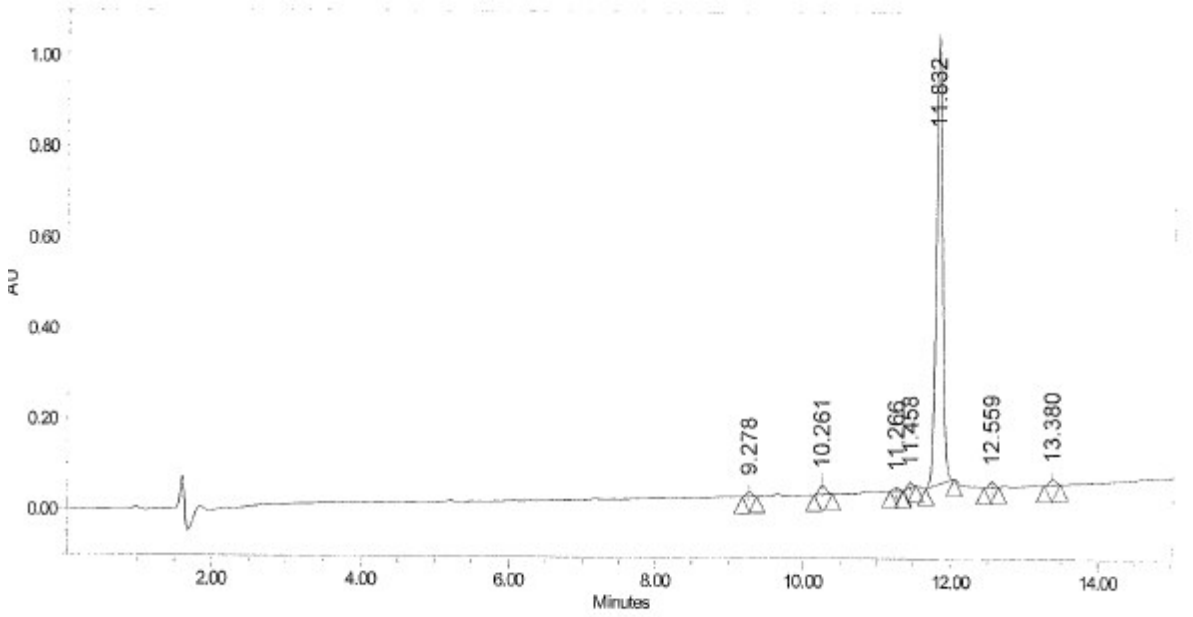


Figure S5. ESI mass spectrum of maleimido propyl $(PEG)_{24}$ dolastatin 10 drug linker with theoretical m/z of MH+ = 2050.18 (please note that in addition to protonated molecular ion peak, both sodium and potassium ion complexed molecular ions peaks were observed.).



Peak Results

	Name	RT	Area	Height	Amount	Units	% Area
1		9.278	57302	10720			0.91
2		10.261	132104	20282			2.10
3		11.266	30717	6417			0.49
4		11.458	52297	10849			0.83
5		11.832	5878353	975585			93.60
6		12.559	61799	11831			0.98
7		13.380	67532	12101			1.08

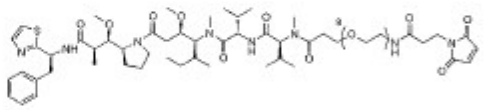


Figure **S6**. HPLC trace of PEG8 Dol10 drug linker.