

Figure S1. ESI-MS spectrum and analytical RP-HPLC spectrum of LCP-1.

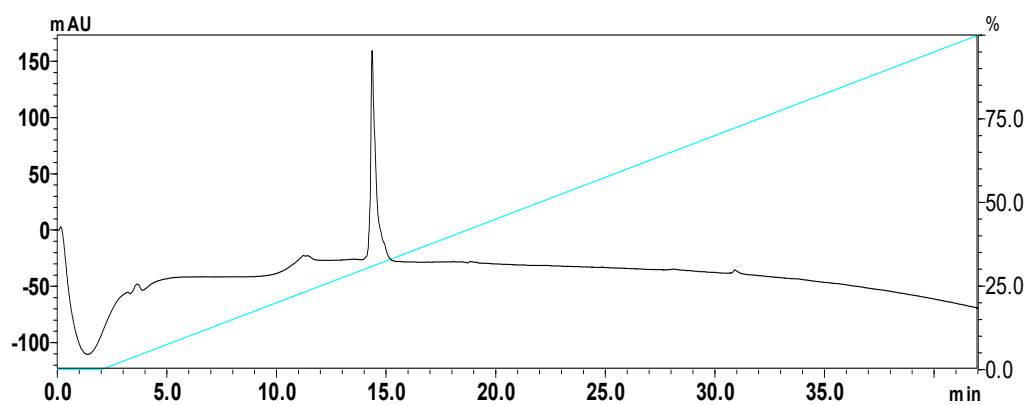
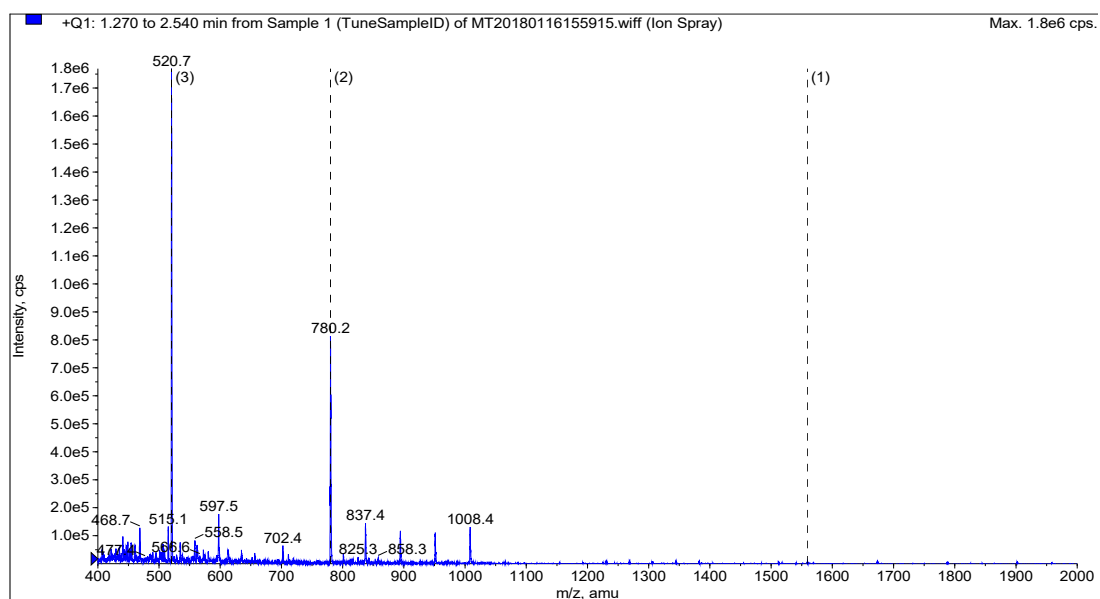


Figure S2. ESI-MS spectrum and analytical RP-HPLC spectrum of Tat₄₇₋₅₇.

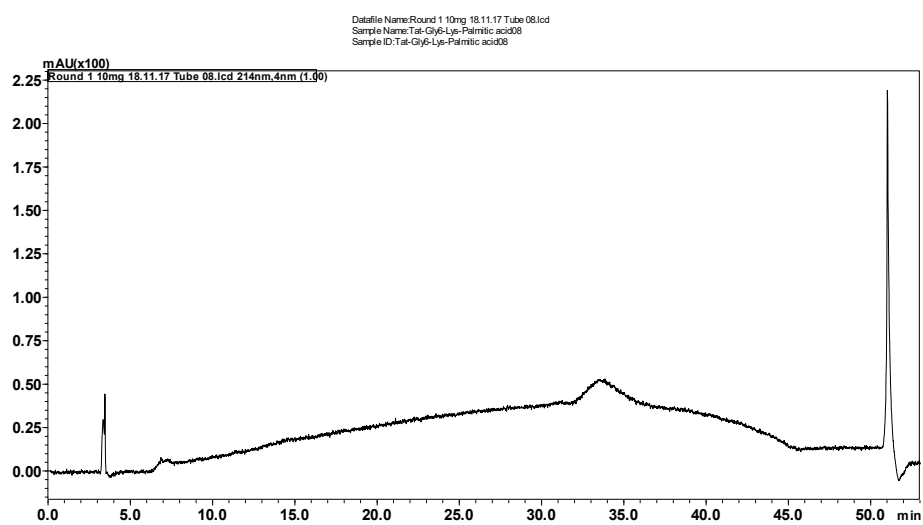
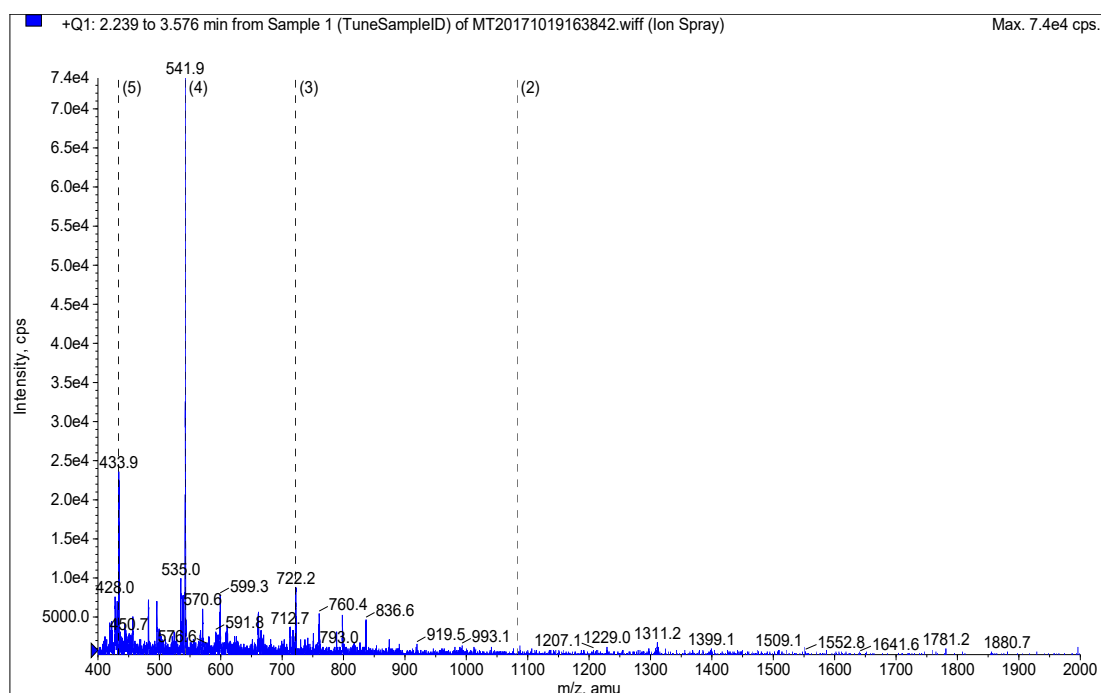


Figure S3. ESI-MS spectrum and analytical RP-HPLC spectrum of compound 1.

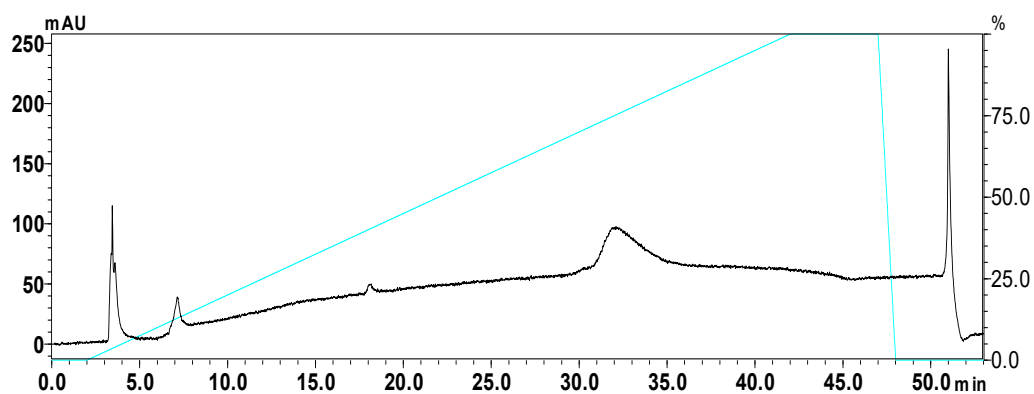
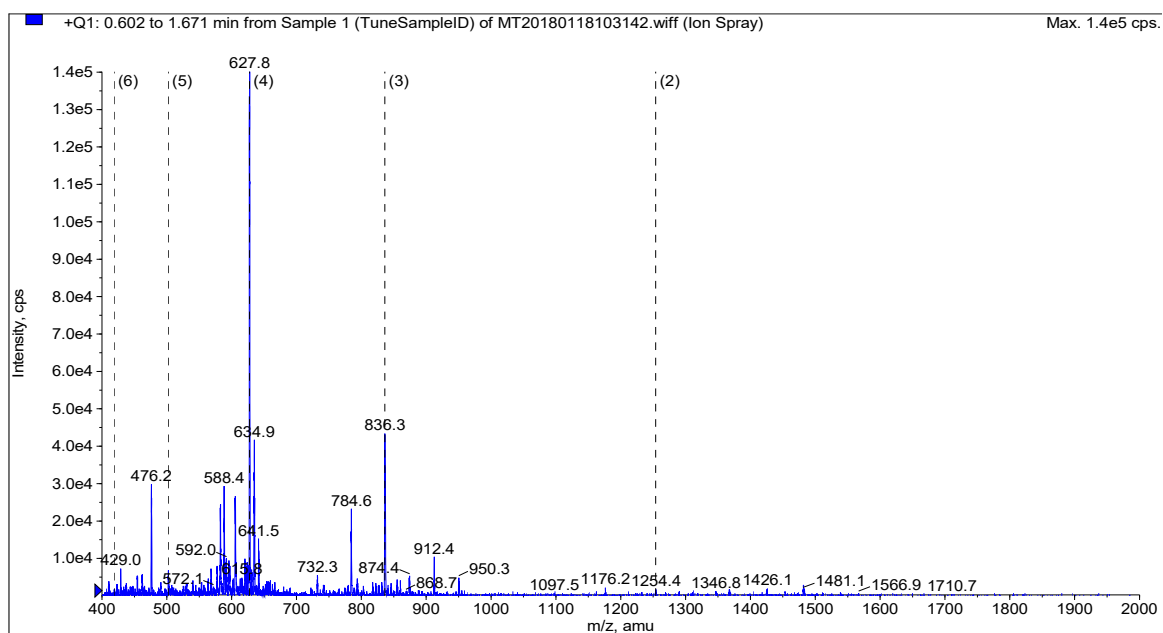


Figure S4. ESI-MS spectrum and analytical RP-HPLC spectrum of compound 2.

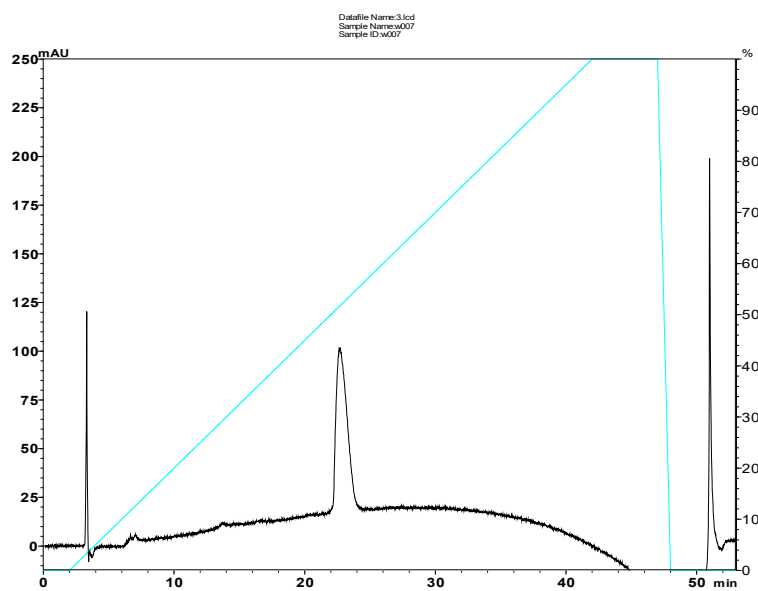
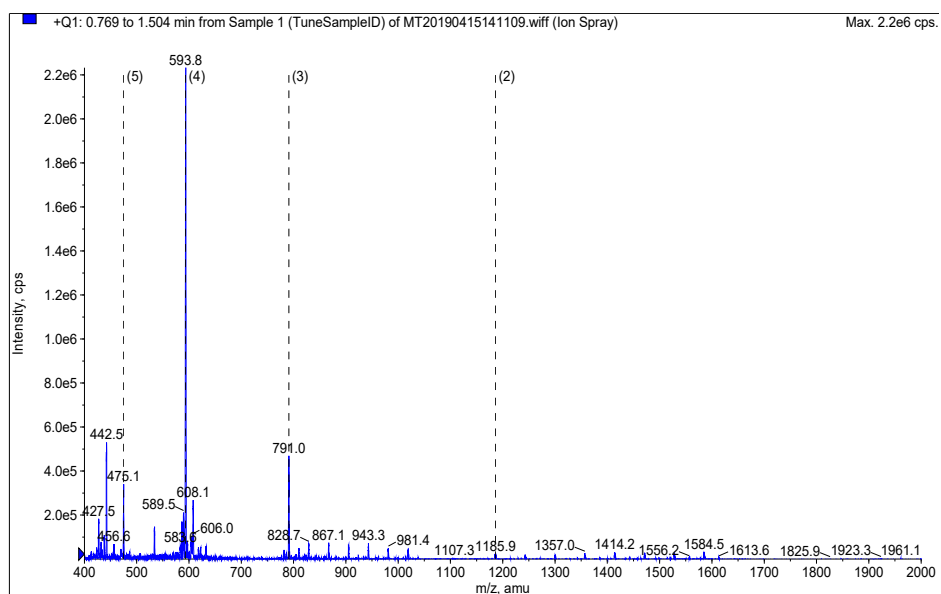


Figure S6. ESI-MS spectrum and analytical RP-HPLC spectrum of compound **4**.

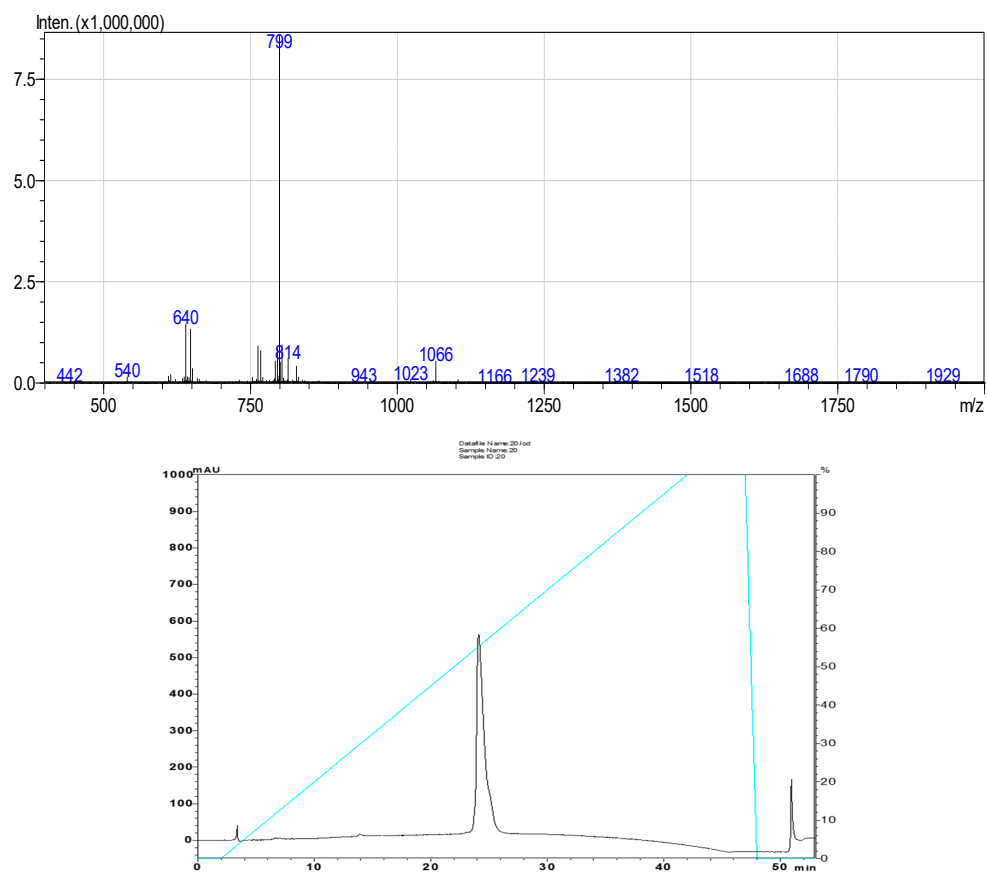


Figure S7. ESI-MS spectrum and analytical RP-HPLC spectrum of compound 5.

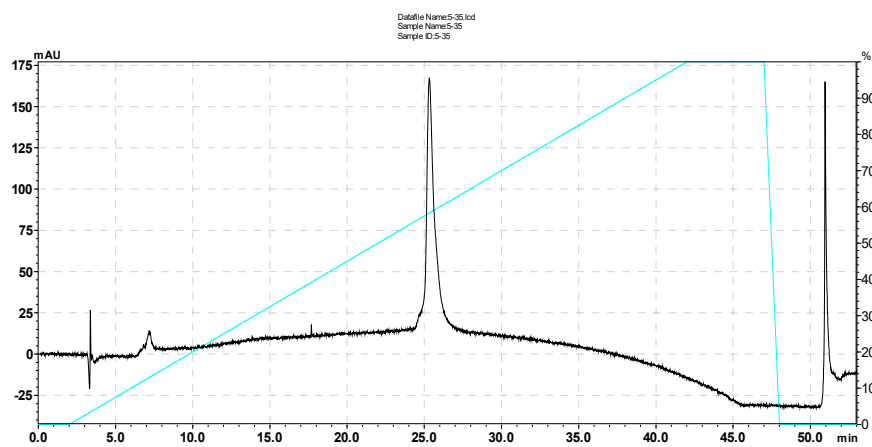
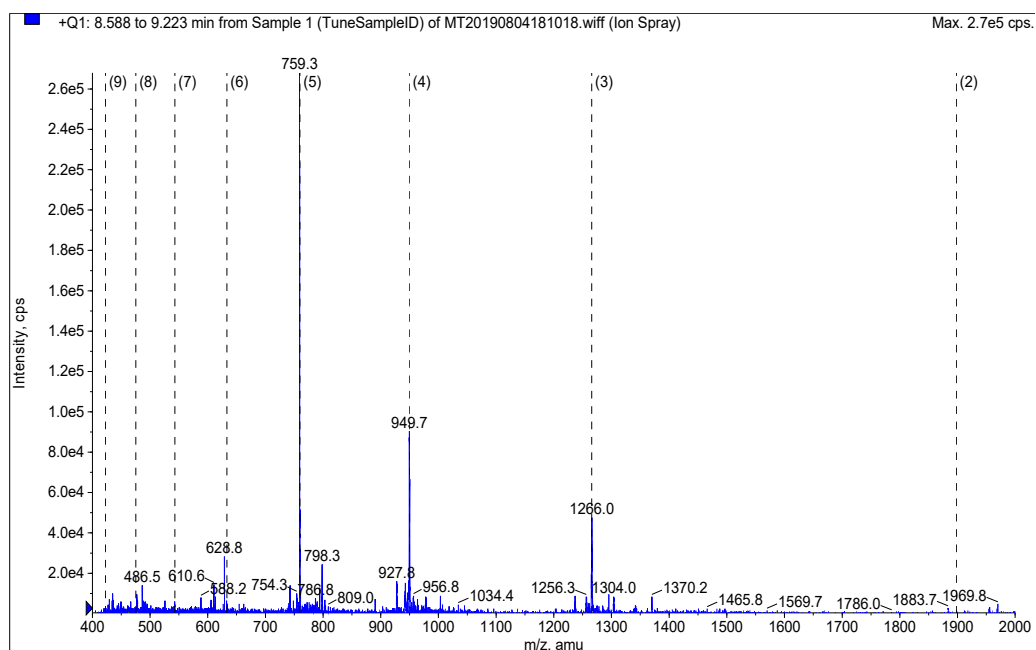


Figure S8. ESI-MS spectrum and analytical RP-HPLC spectrum of compound 6.

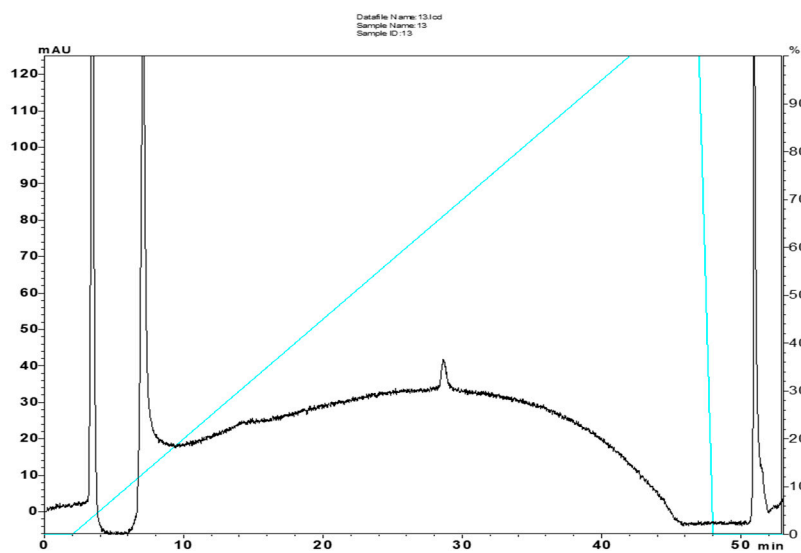
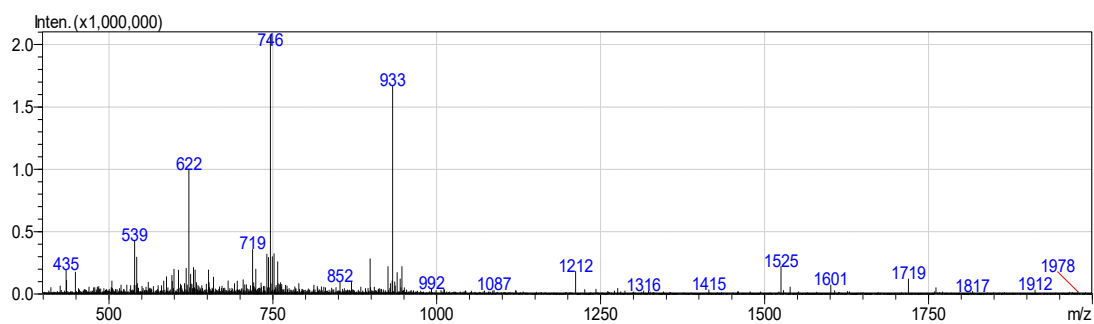
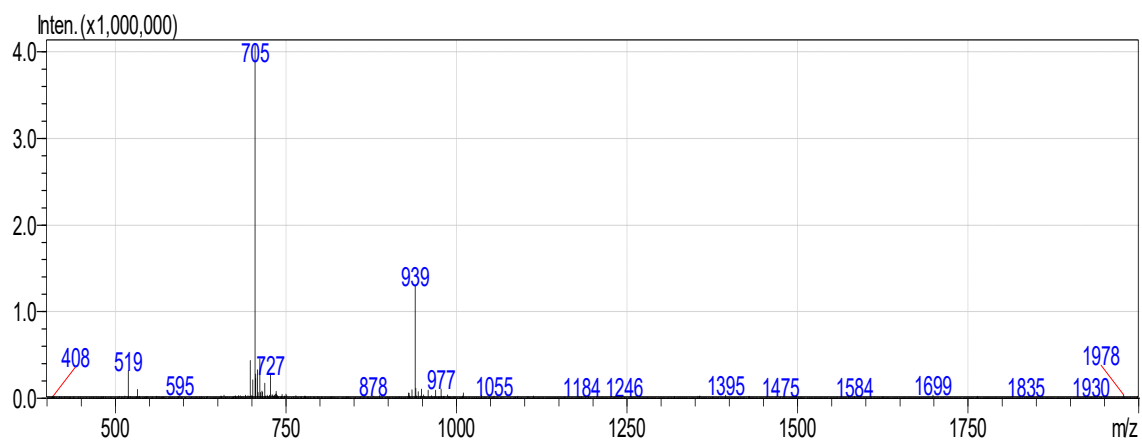


Figure S9. ESI-MS spectrum and analytical RP-HPLC spectrum of compound 7.



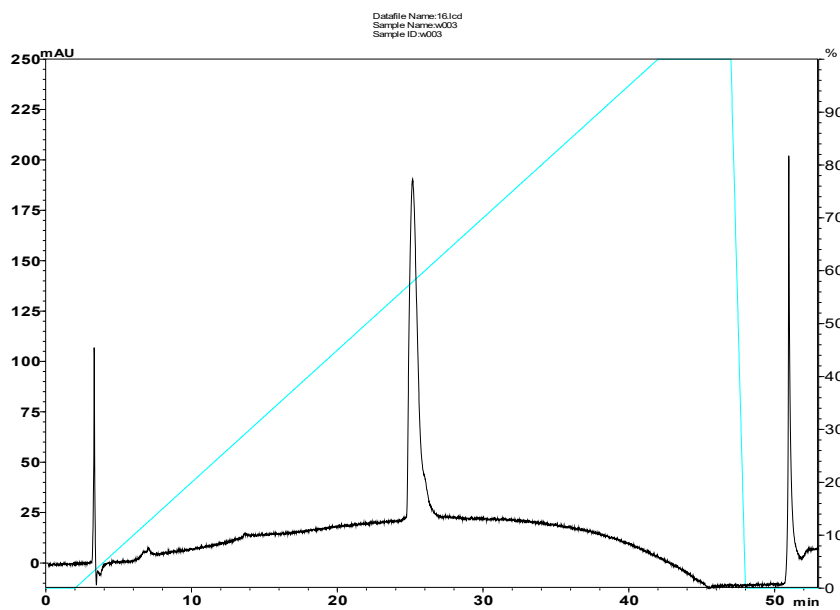


Figure S10. ESI-MS spectrum and analytical RP-HPLC spectrum of compound 8.

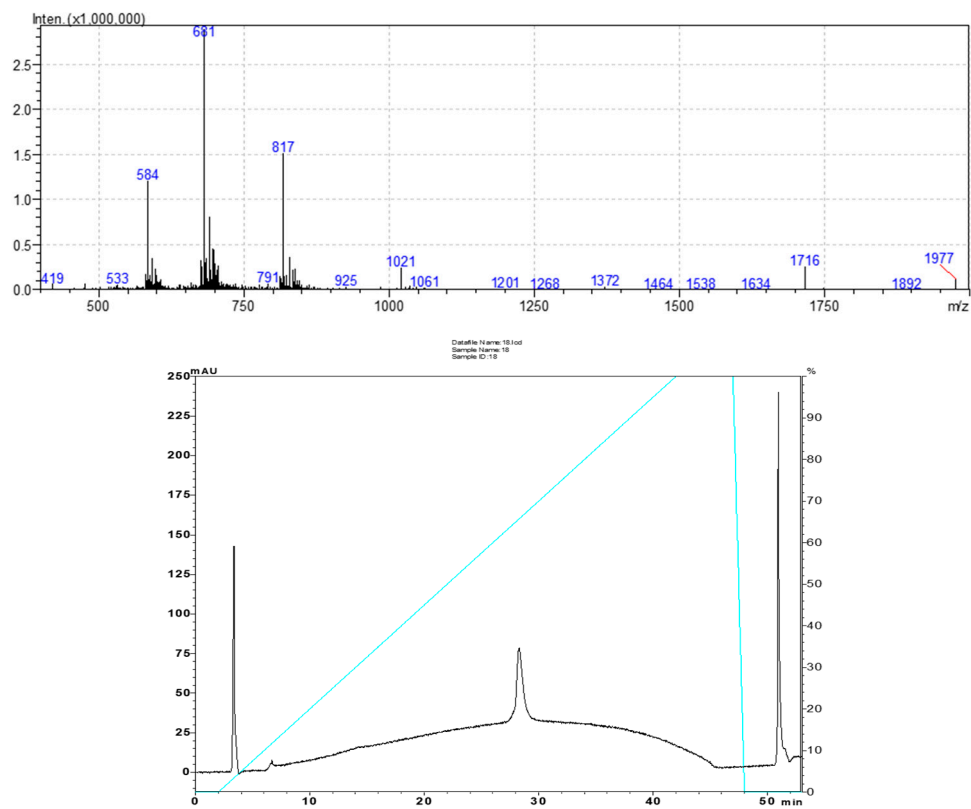
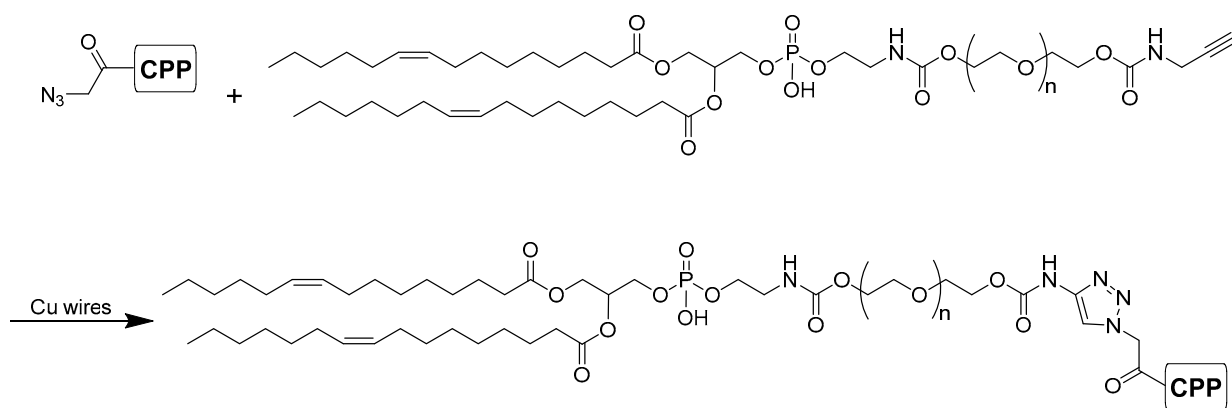


Figure S11. ESI-MS spectrum and analytical RP-HPLC spectrum of compound 9.



10: $n \approx 77$, CPP = Tat₄₇₋₅₇

Figure S12. The click chemistry of compound **10**. Compound **10** was produced based on a copper-catalyzed alkyne-azide 1,3-dipolar cycloaddition (CuAAC) “click” reaction to conjugate DOPE-PEG₃₄₀₀-alkyne to the azide-Tat₄₇₋₅₇.

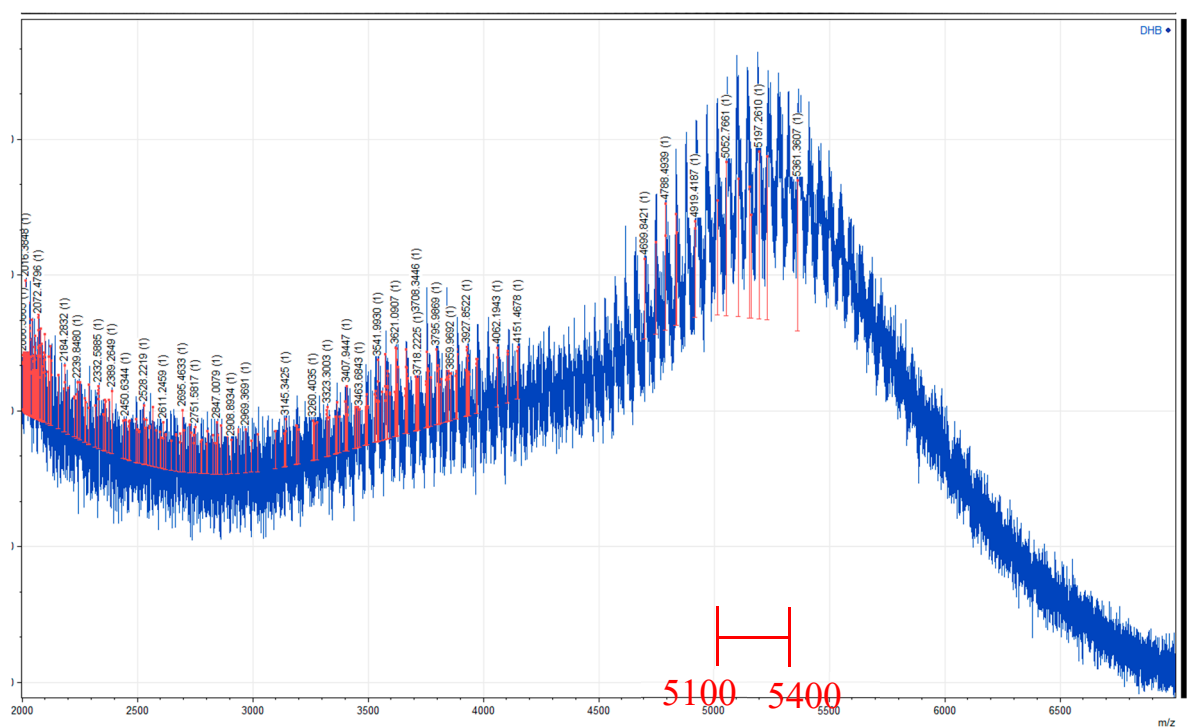


Figure S13. The MALDI-TOF Mass Spectrometry of DOPE-PEG₃₄₀₀-Tat₄₇₋₅₇ (molecular weight Range: 5558-5758) (compound 10)

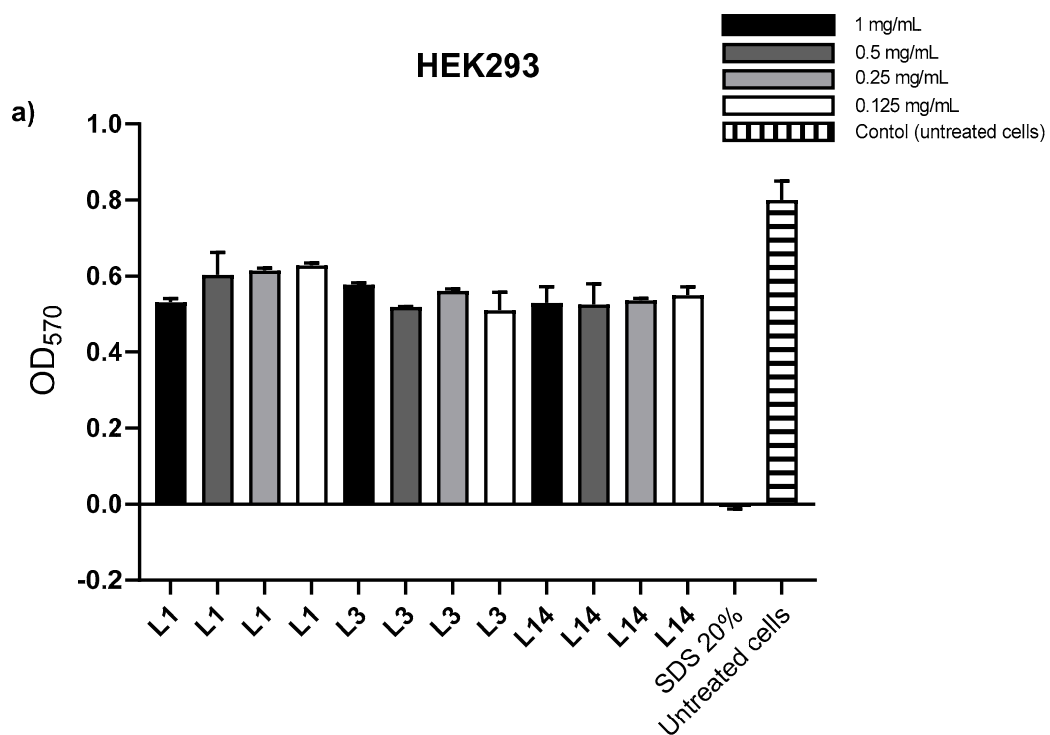


Figure S14. The cytotoxicity results for L1, L3 and L14 against HEK-293: human kidney cell line. Values are reported as mean: standard error of the mean (SEM).

