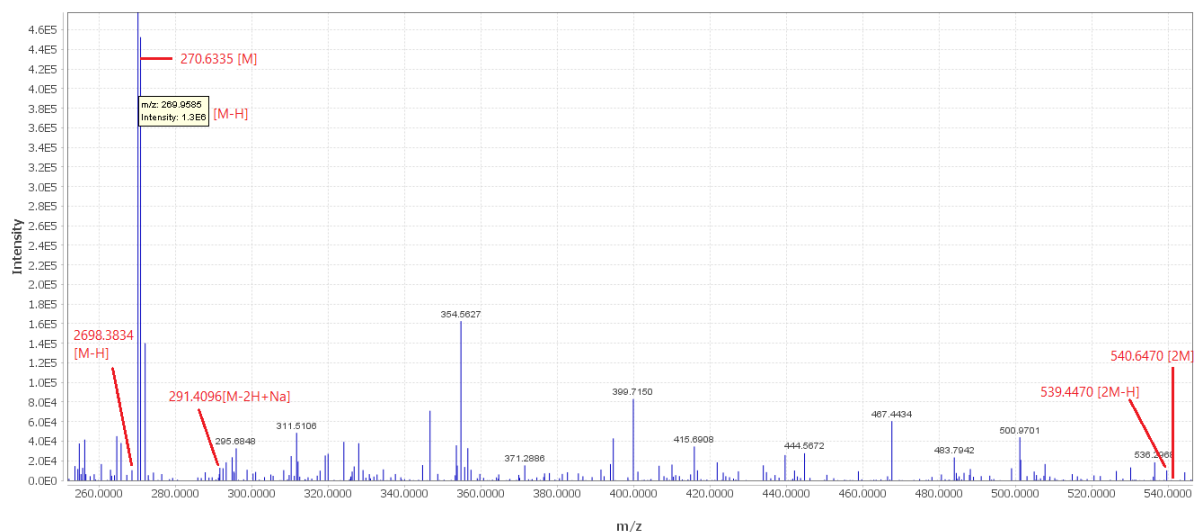
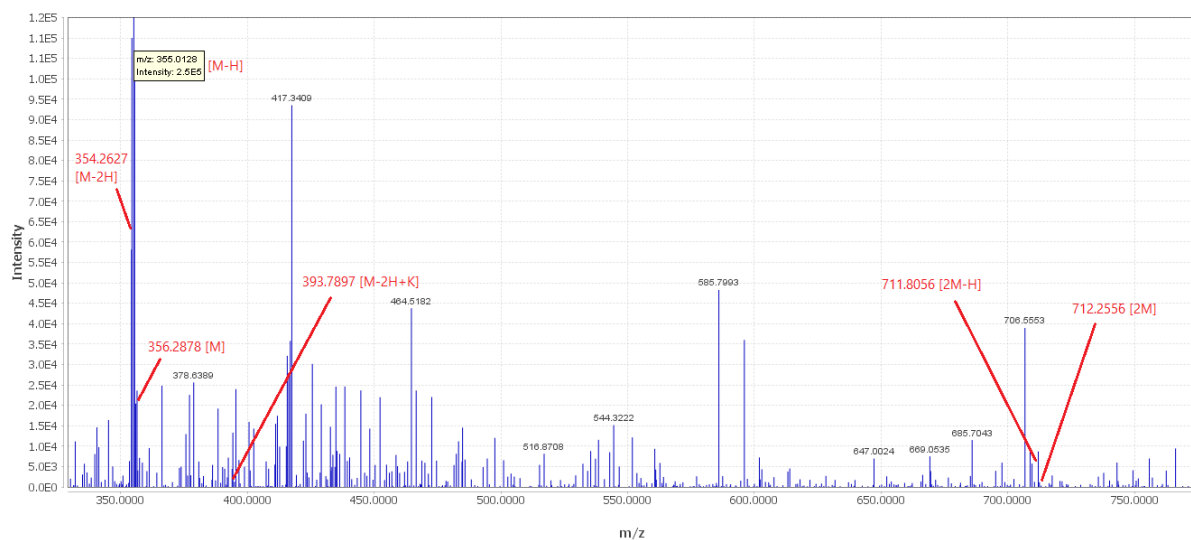
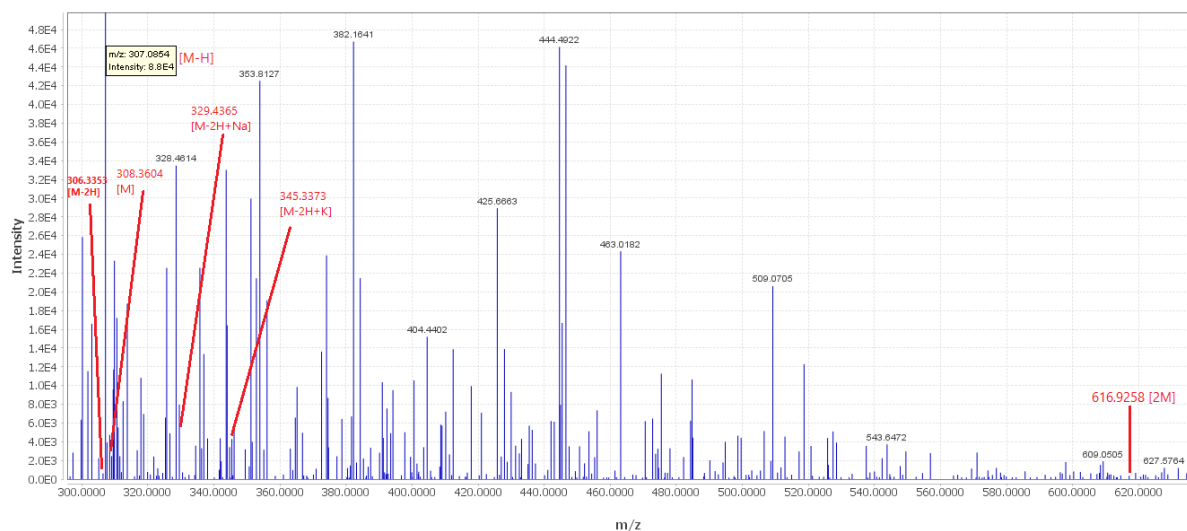
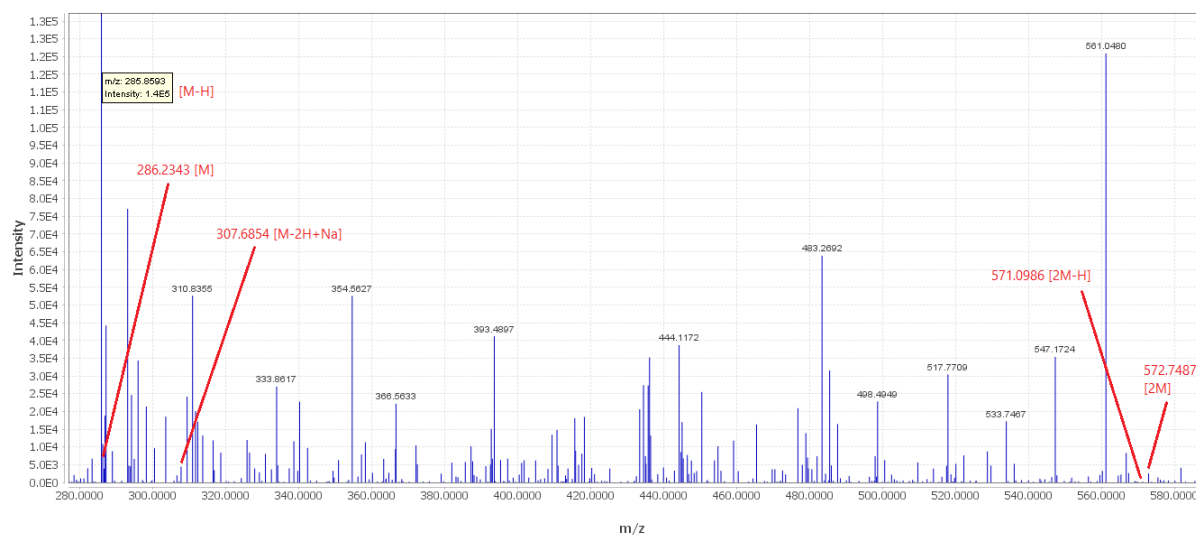
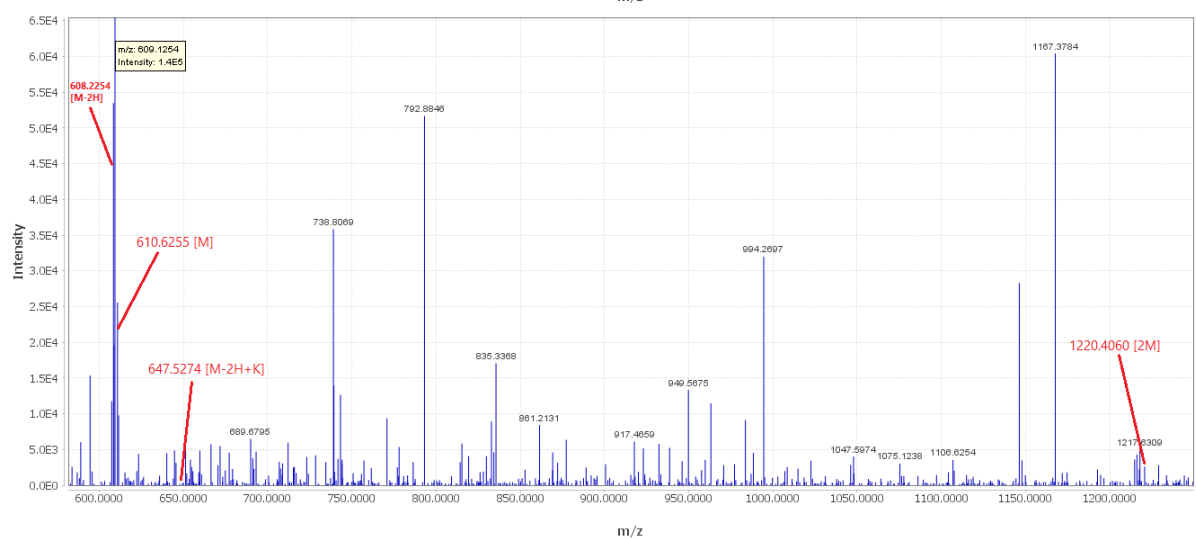
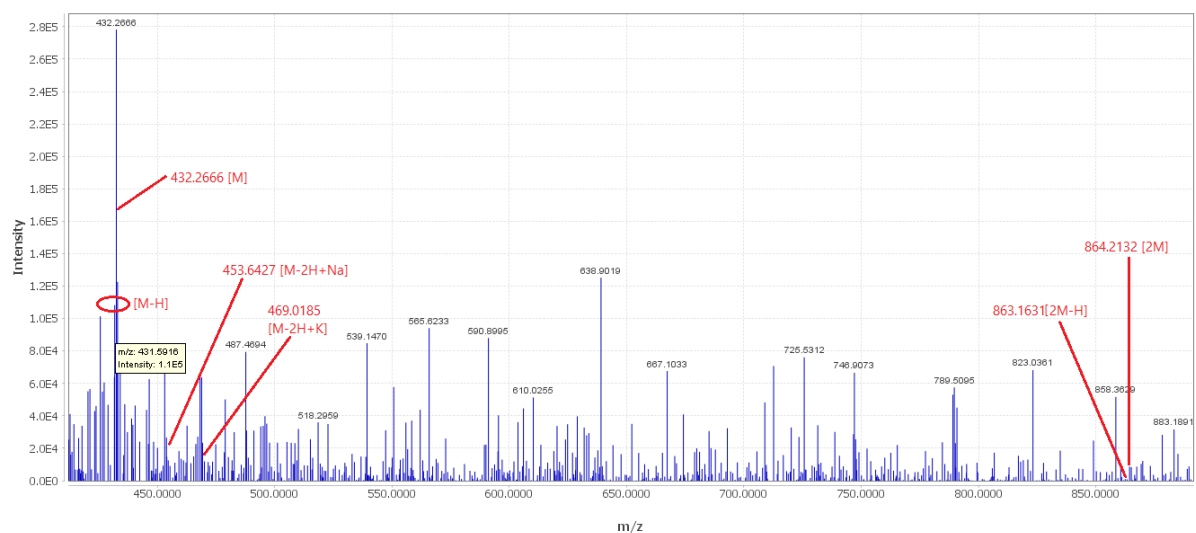
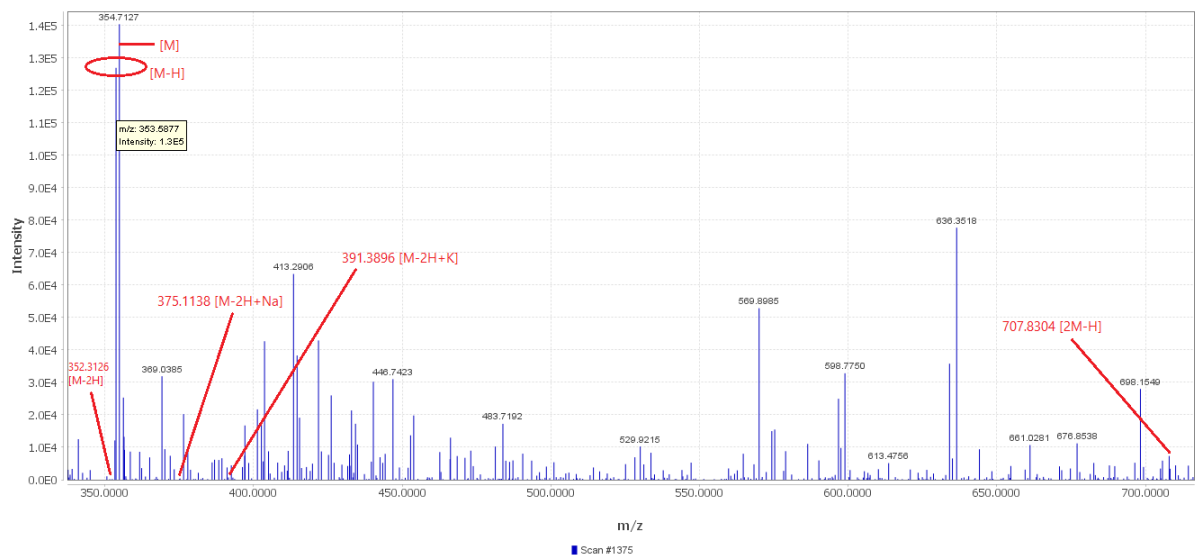
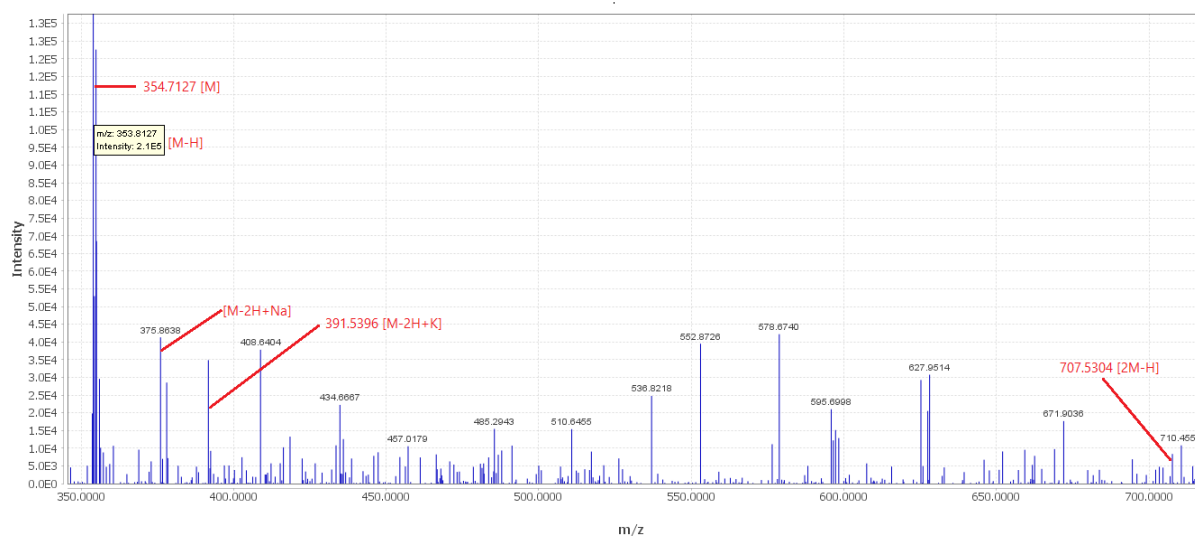
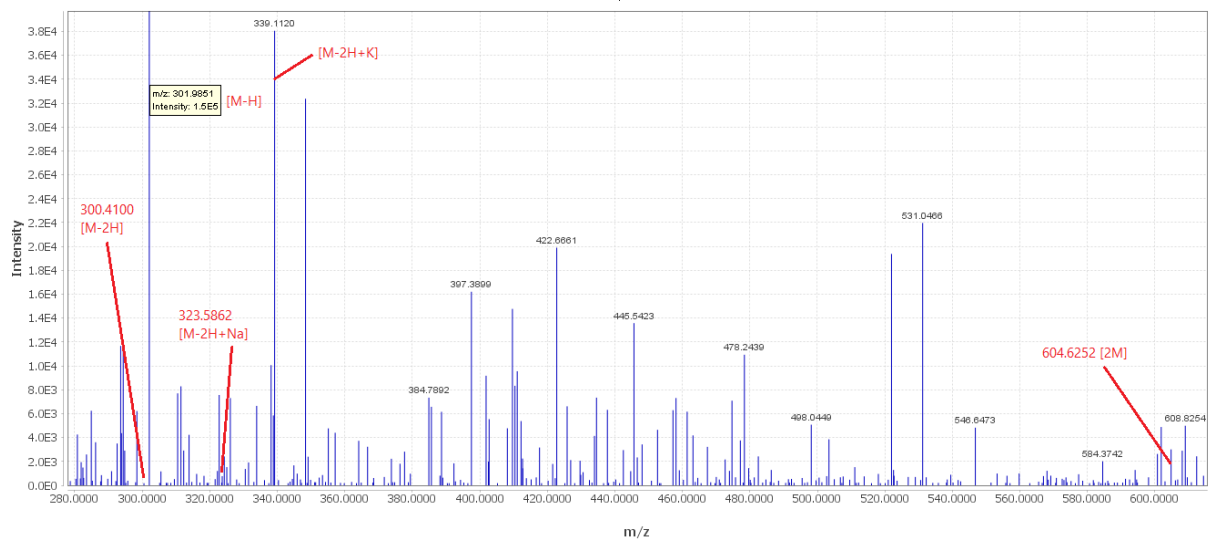


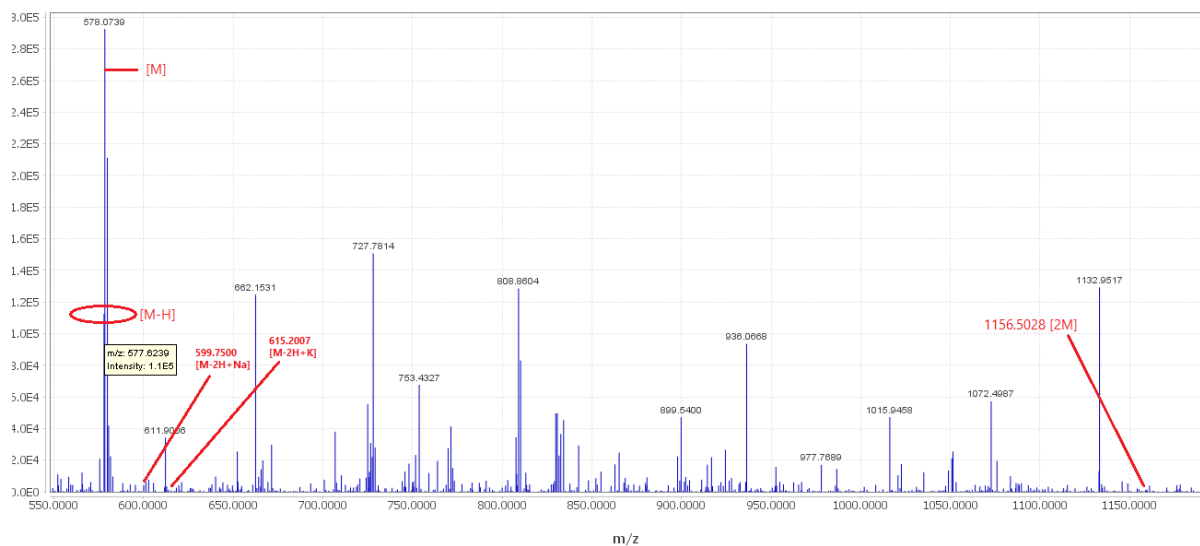
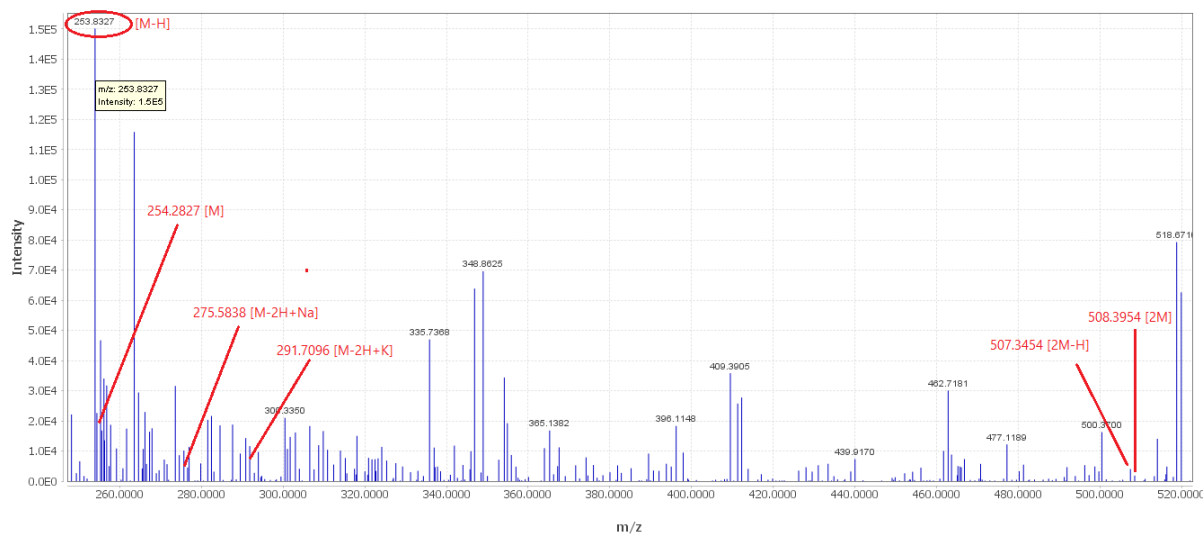
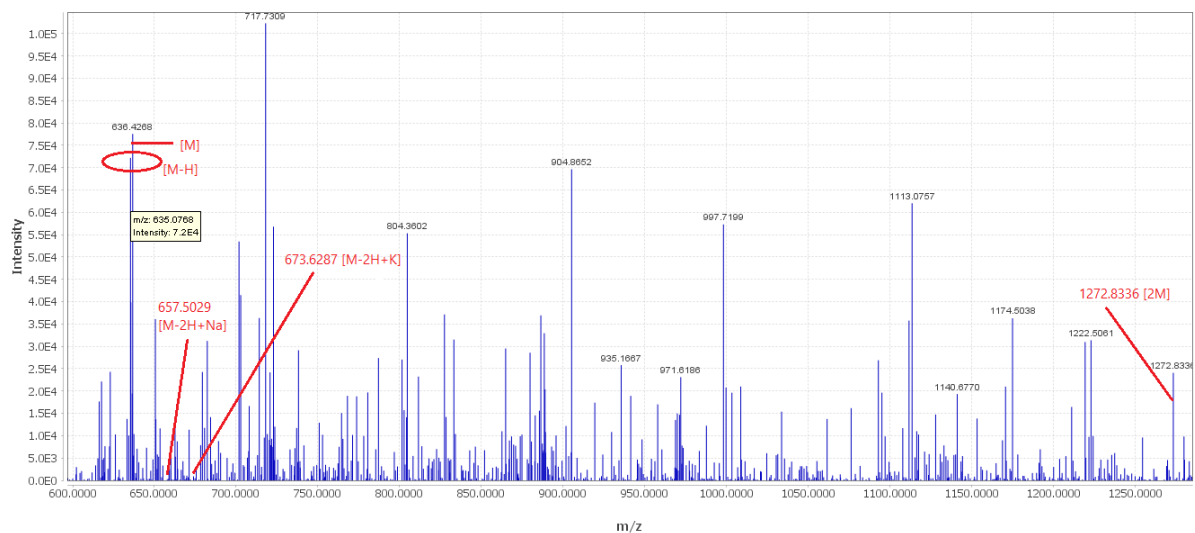
Table S1- Phytochemical compounds identified in the methanolic extract of *Silene conoidea* root.

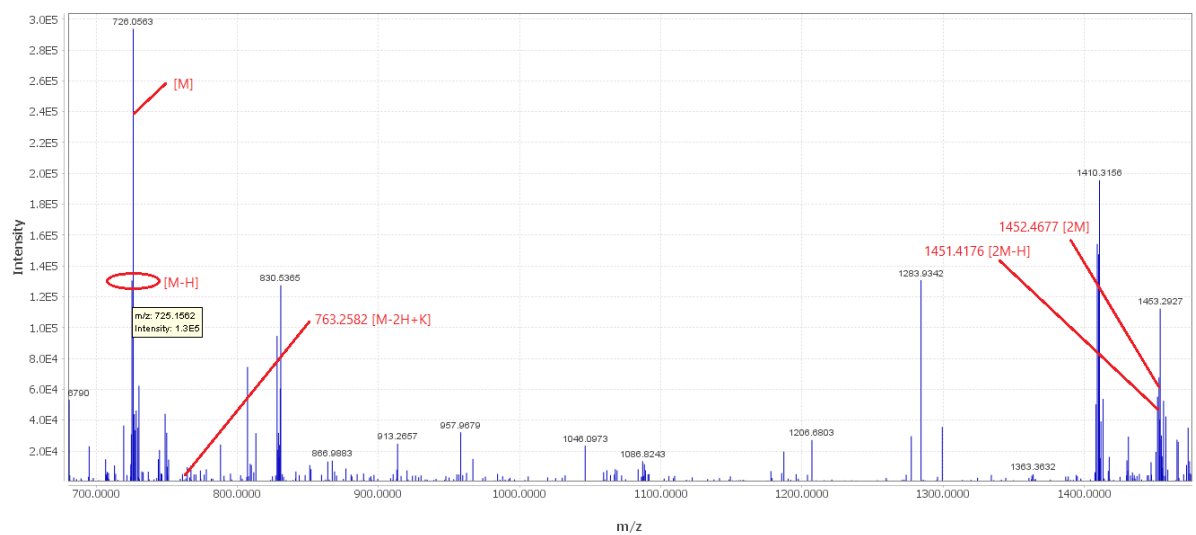
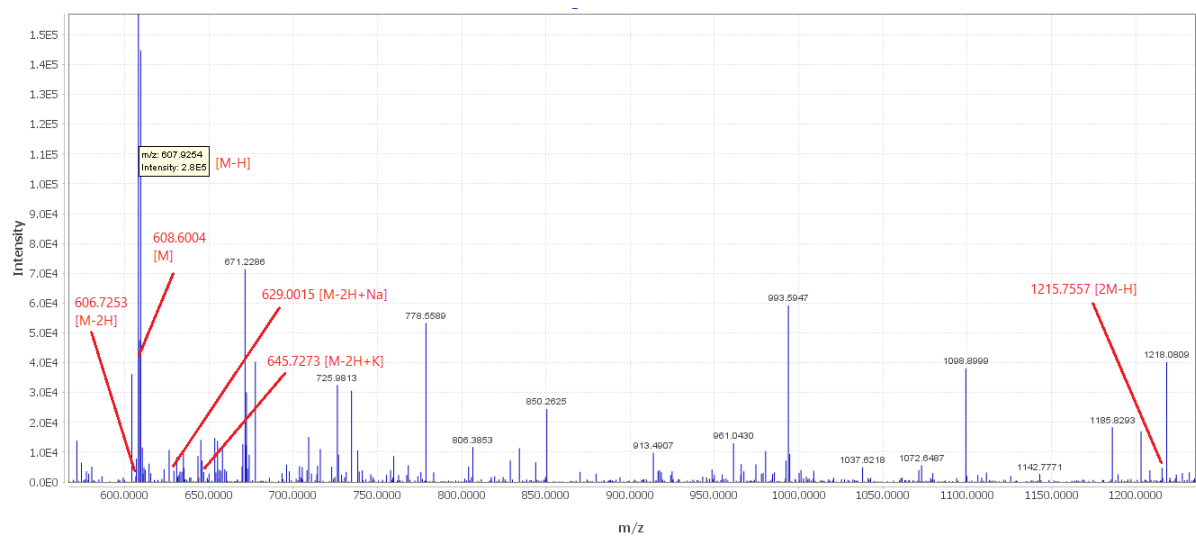
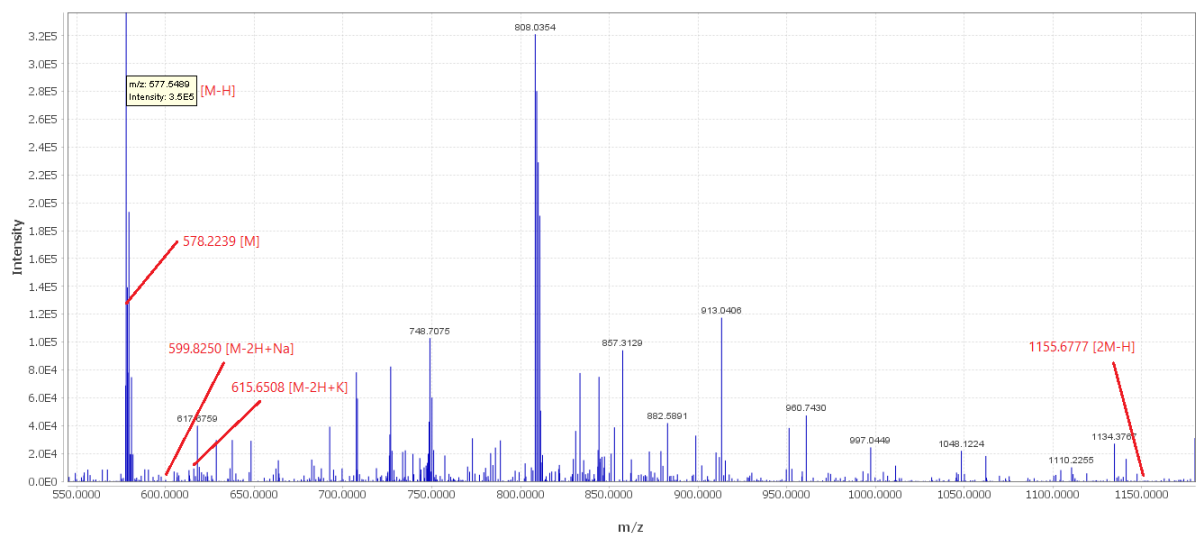
Chlorogenic acid	Vanillic acid hexoside	Esculin	Rutin	Diosmin
Luteolin-di-O glucoside derivative	Orientin-O-glucoside derivative	Isovitexin-O-glucoside	Isosaponarin	Vitexin-O-glucoside
Vitexin-O-rhamnoside derivative	Isovitexin-O-rhamnoside derivative	Vicenin derivative	Swertisin-O-glucoside derivative	Apigenin-O-glucoside
Vitexin	Isovitexin	Neovitexin	Isonovitexin	Luteolin
Hesperidin	Hesperetin	Apigenin	Procyanidin B1	Sileneoside H
Sileneoside E	Sileneoside C	Sileneoside B	20-Hydroxyecdysone-25- Glucoside	Sileneoside F
Gypsogenin-O-glucuronide derivative	Sinocrassulose X			

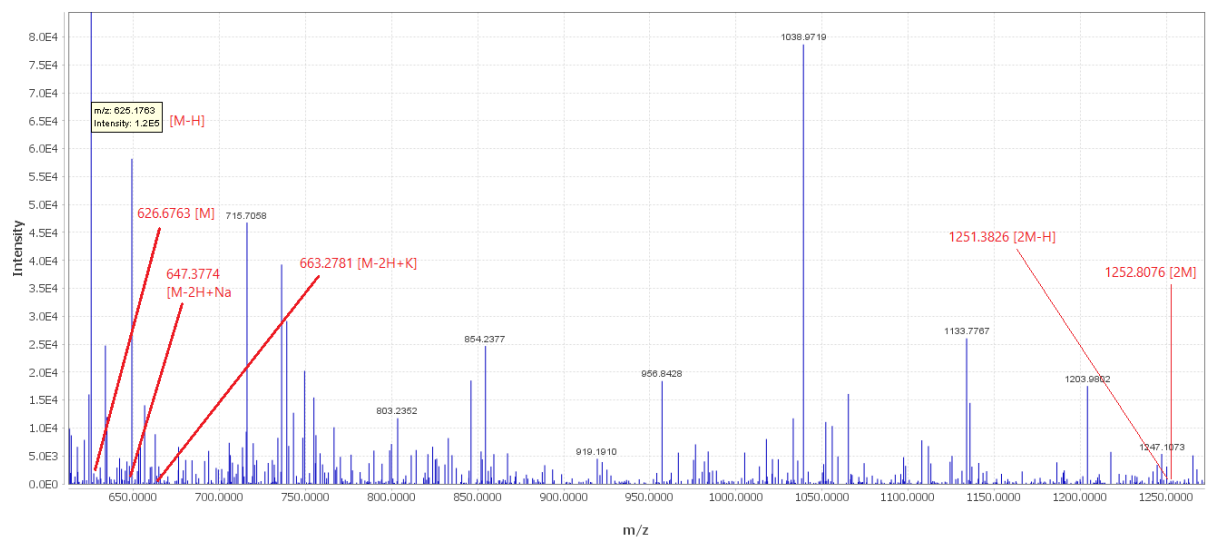
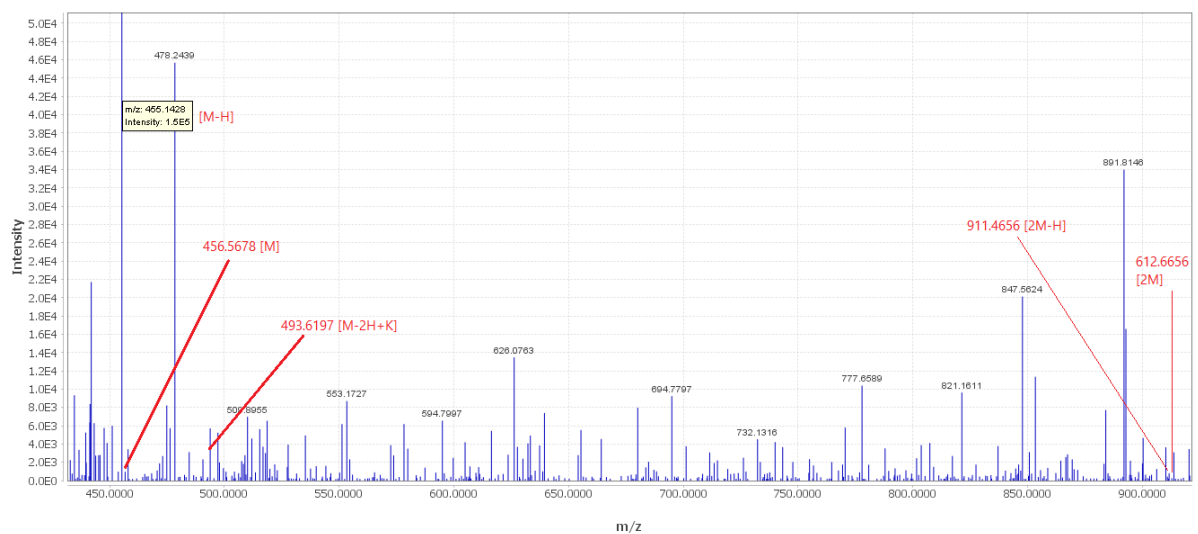
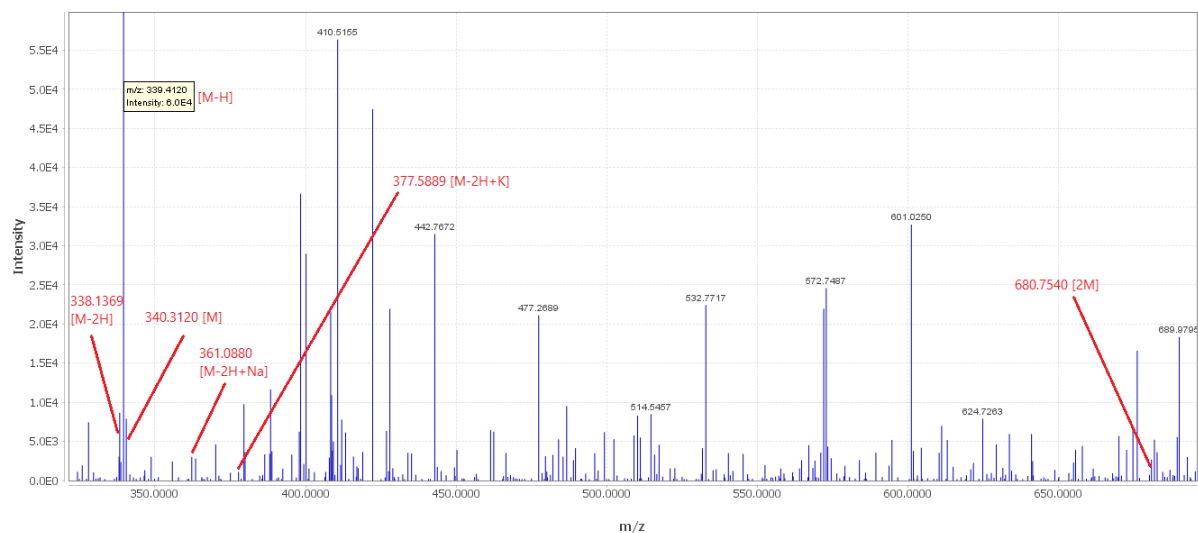


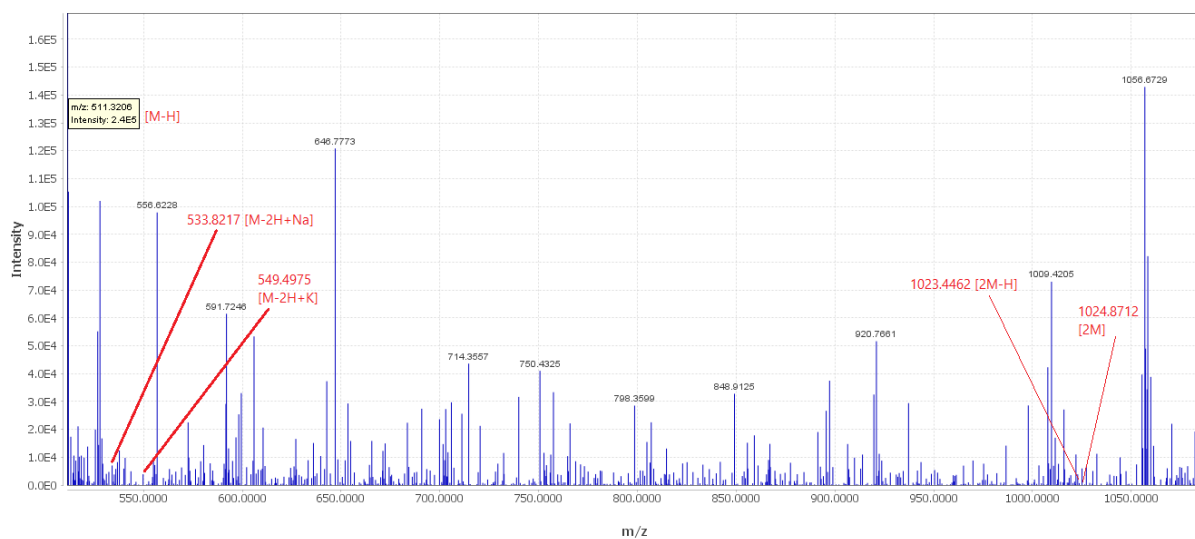
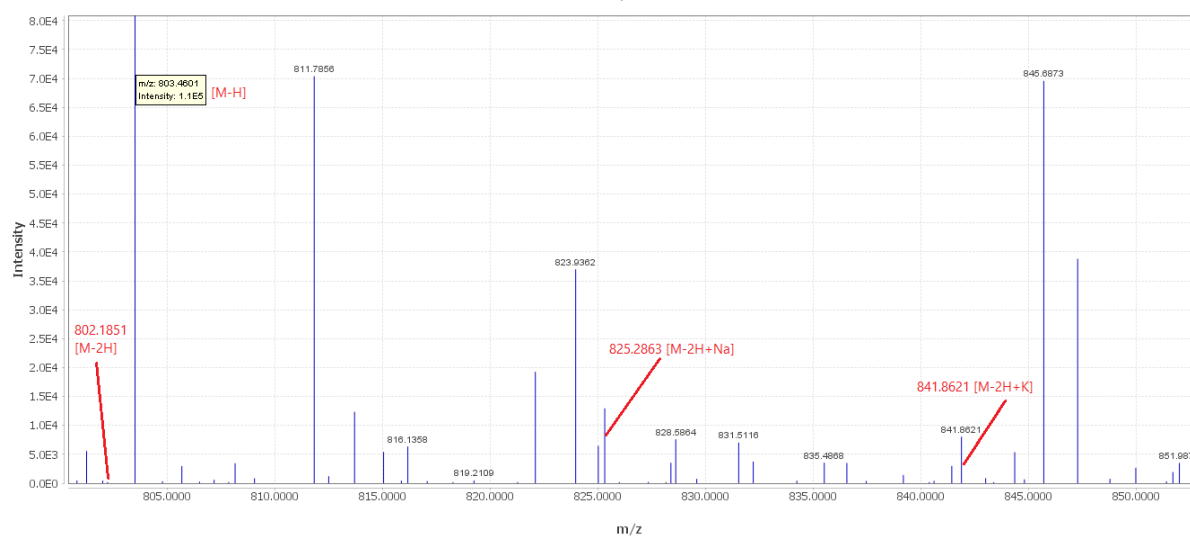
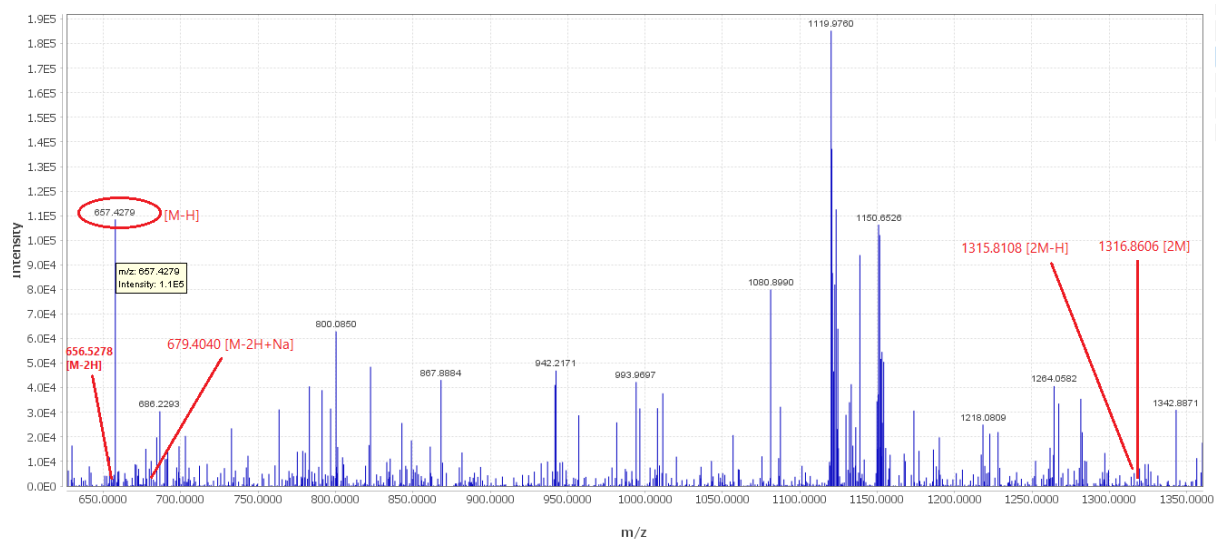


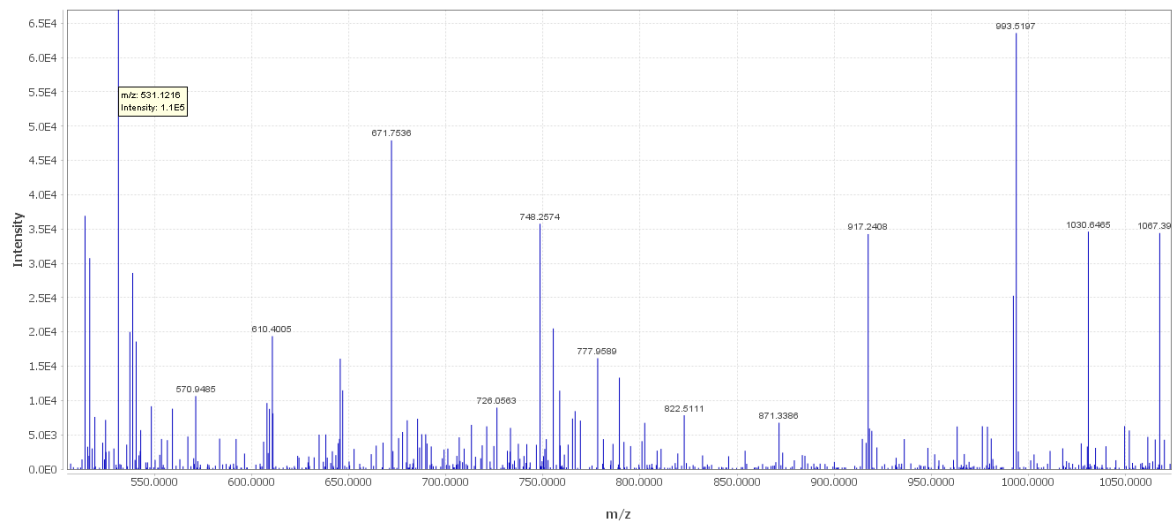
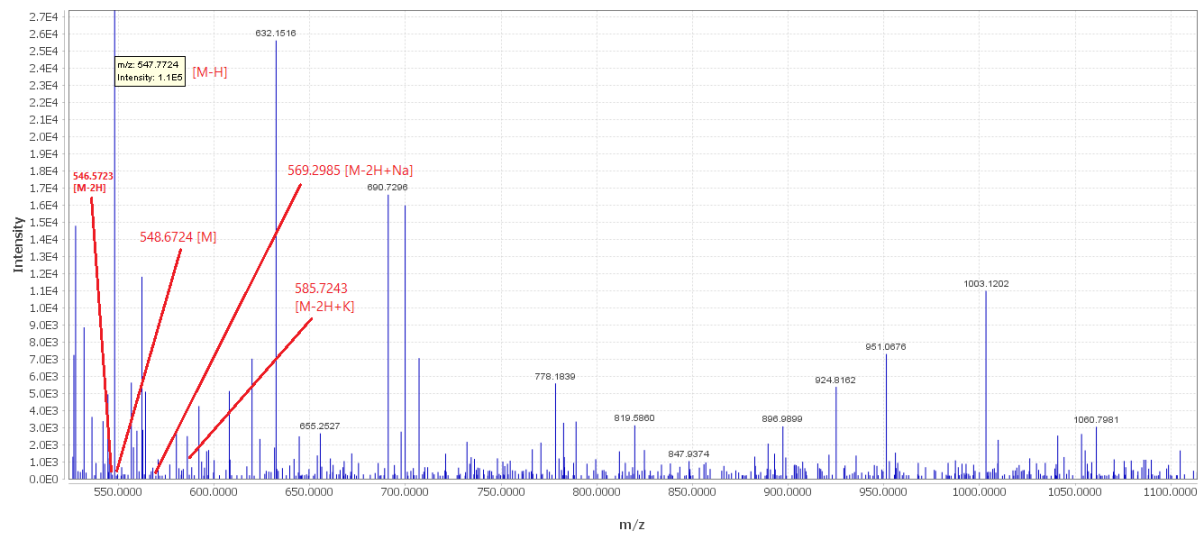
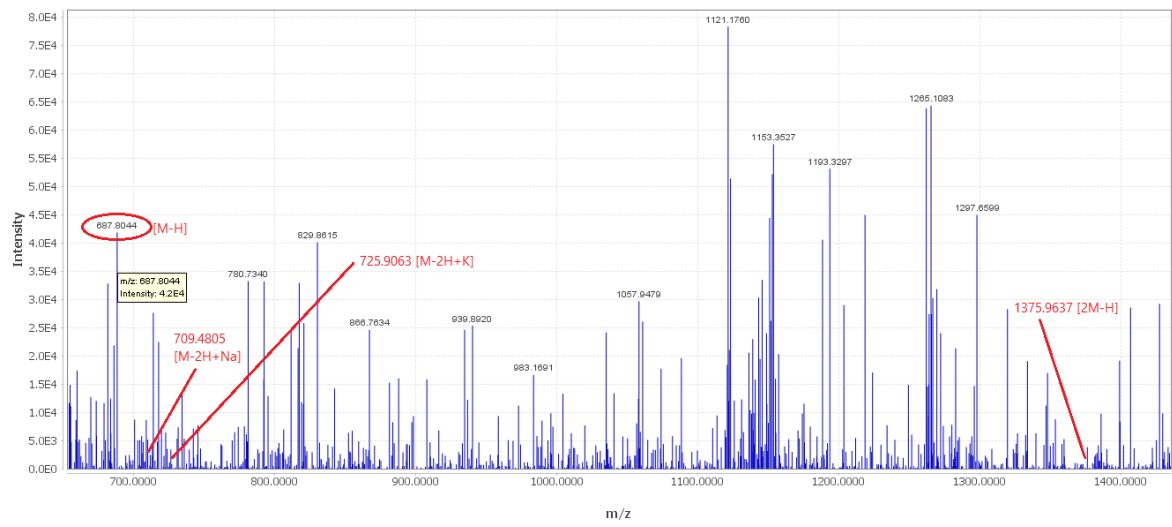


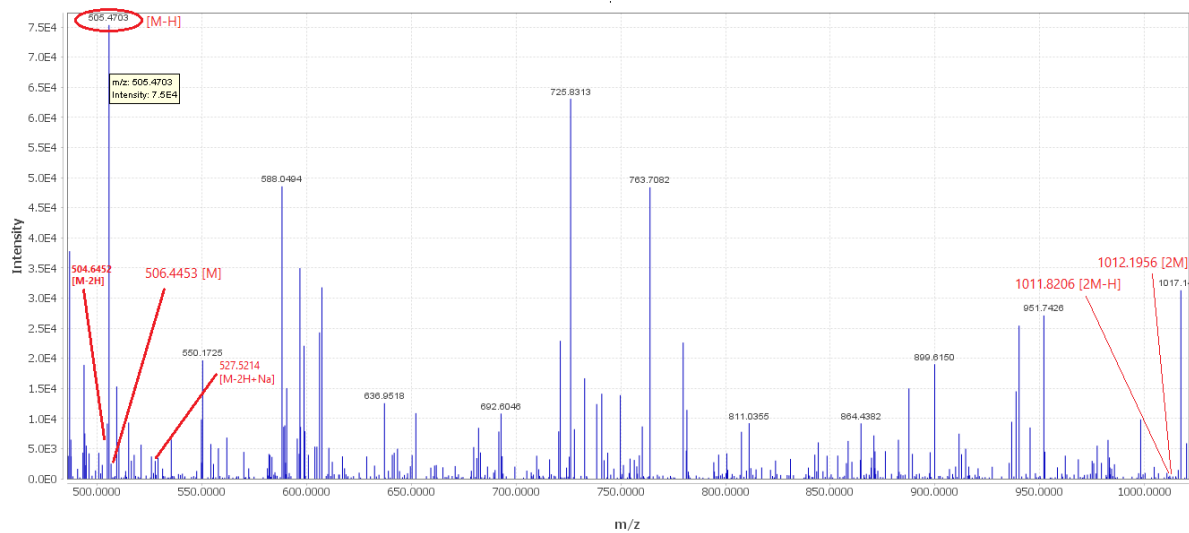
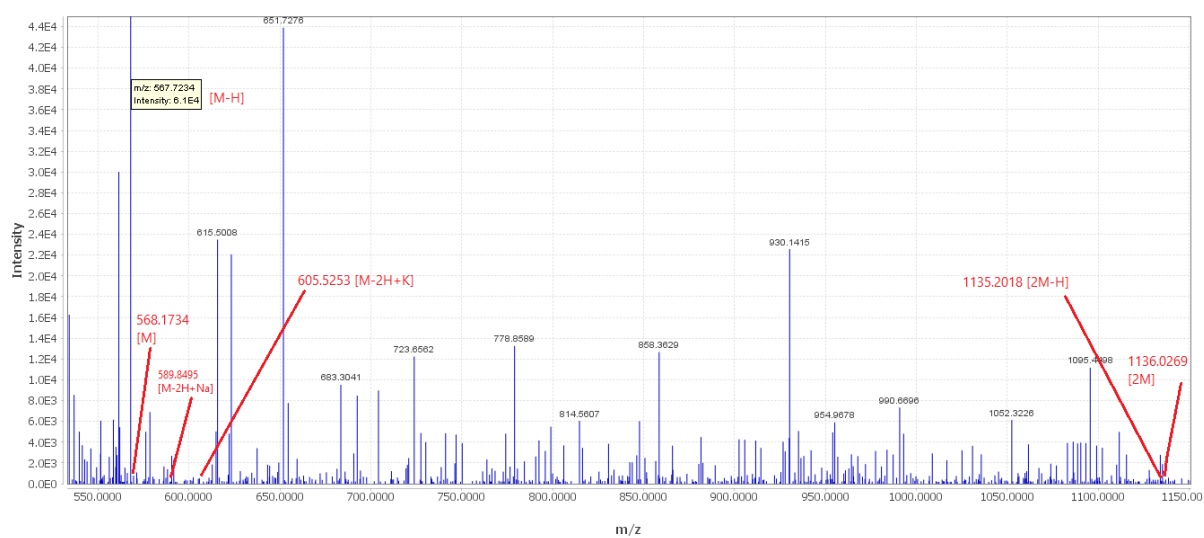
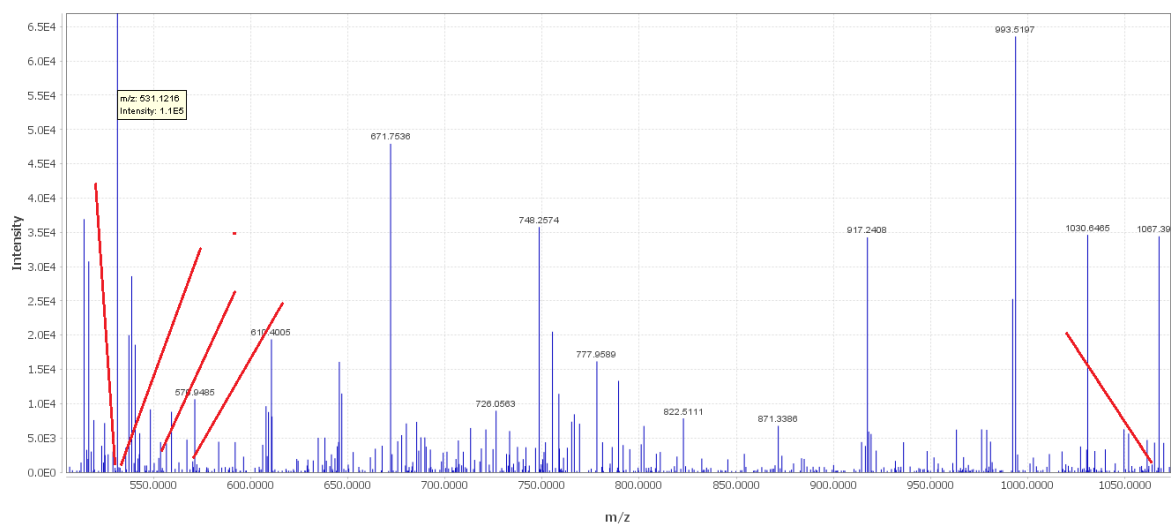


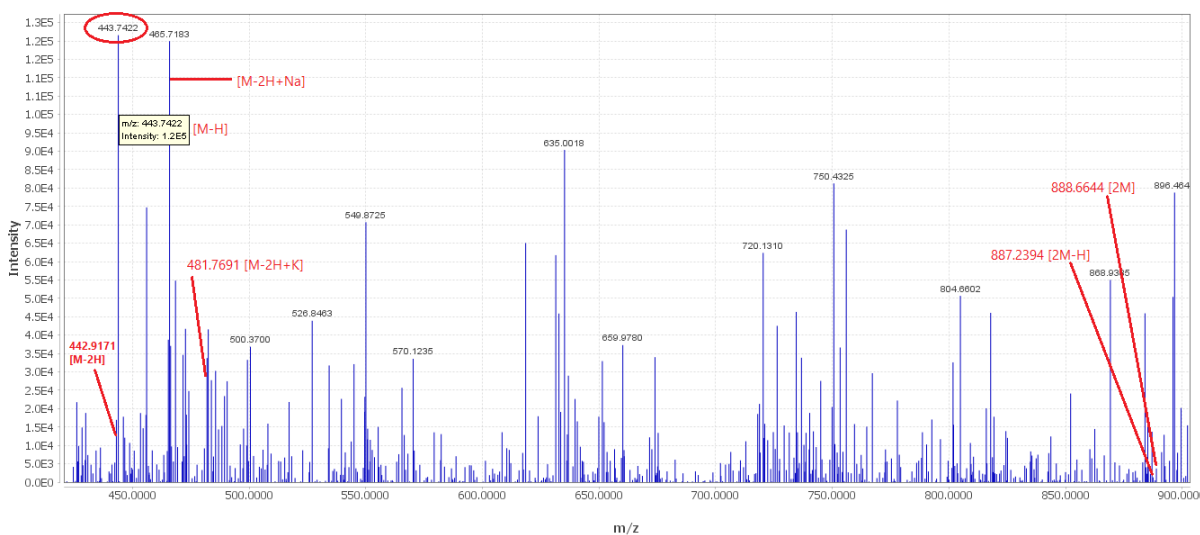
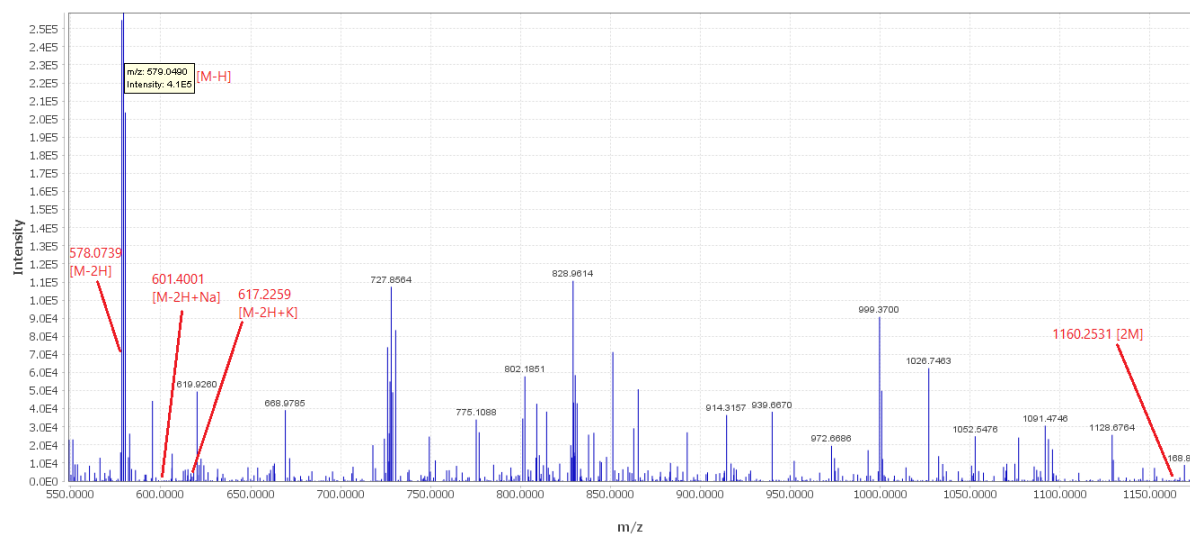
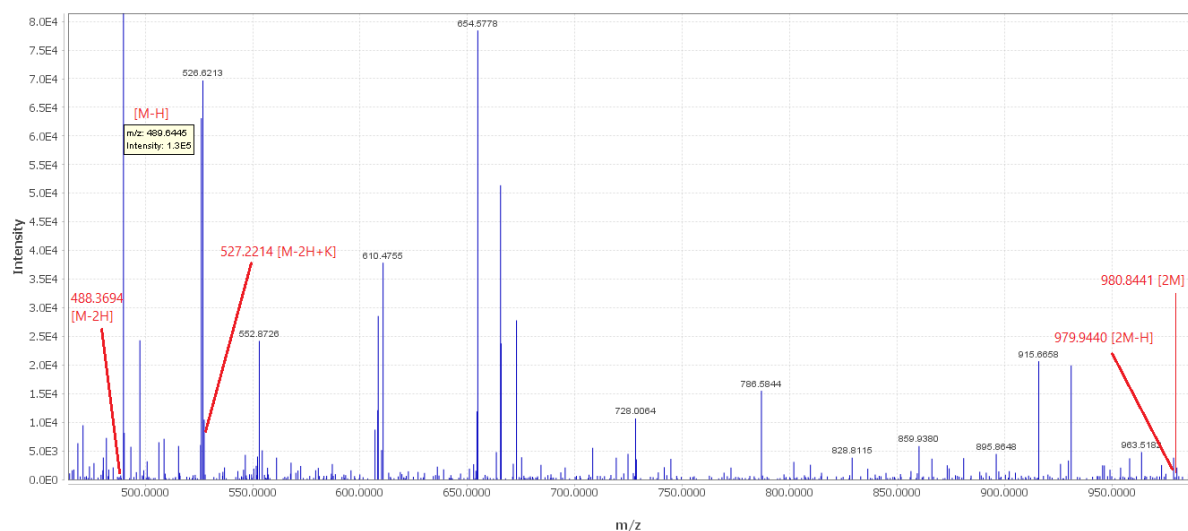


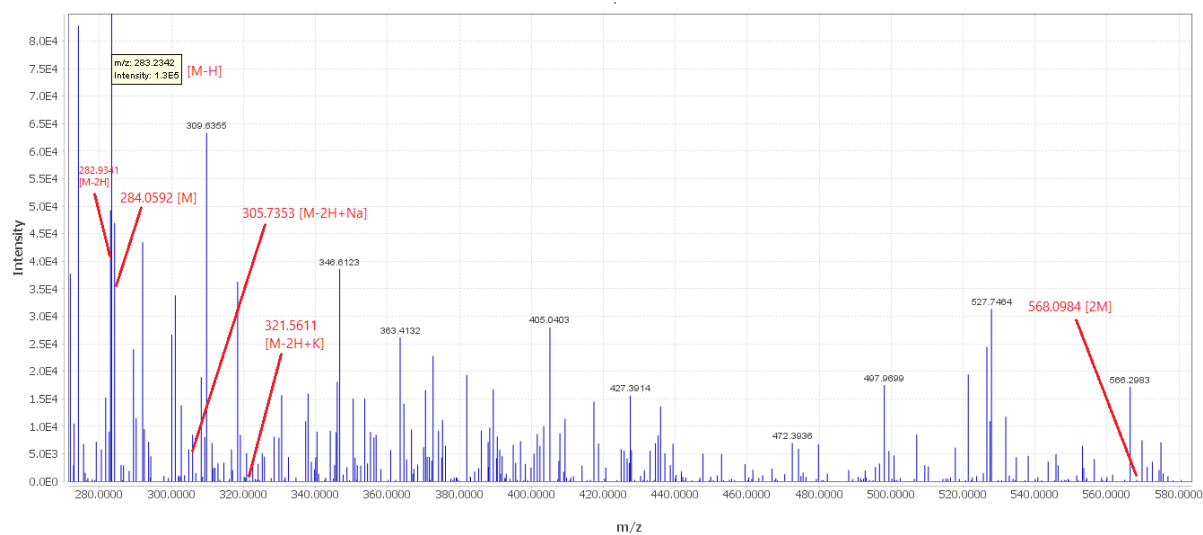
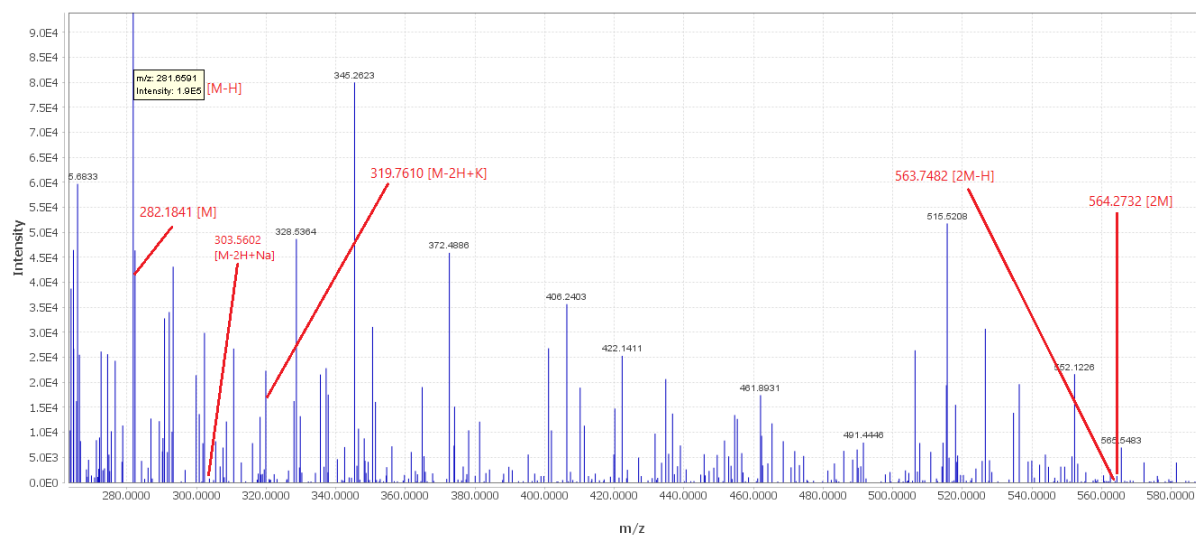
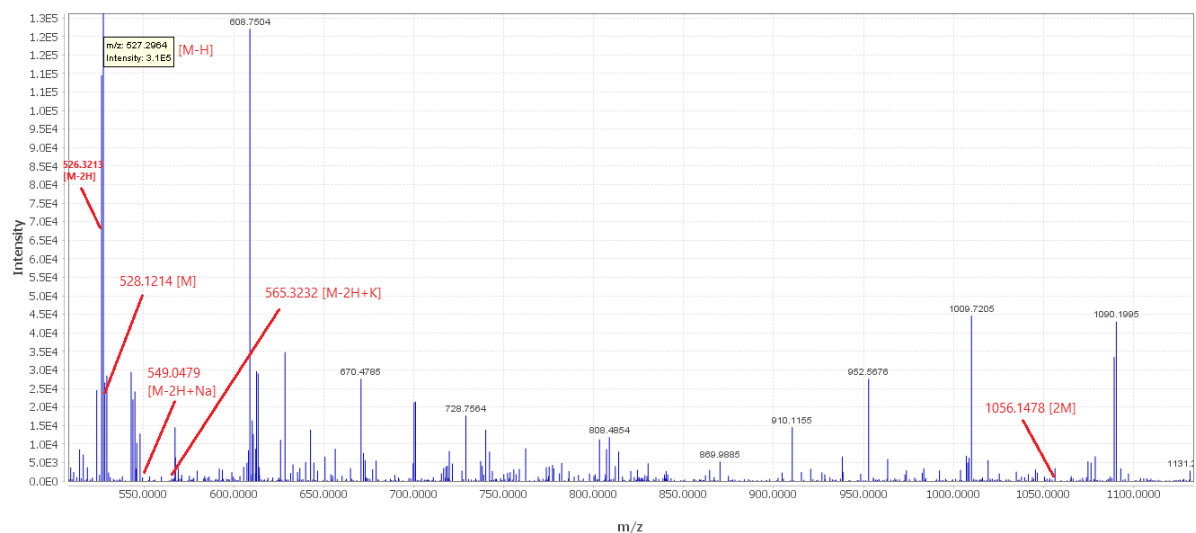


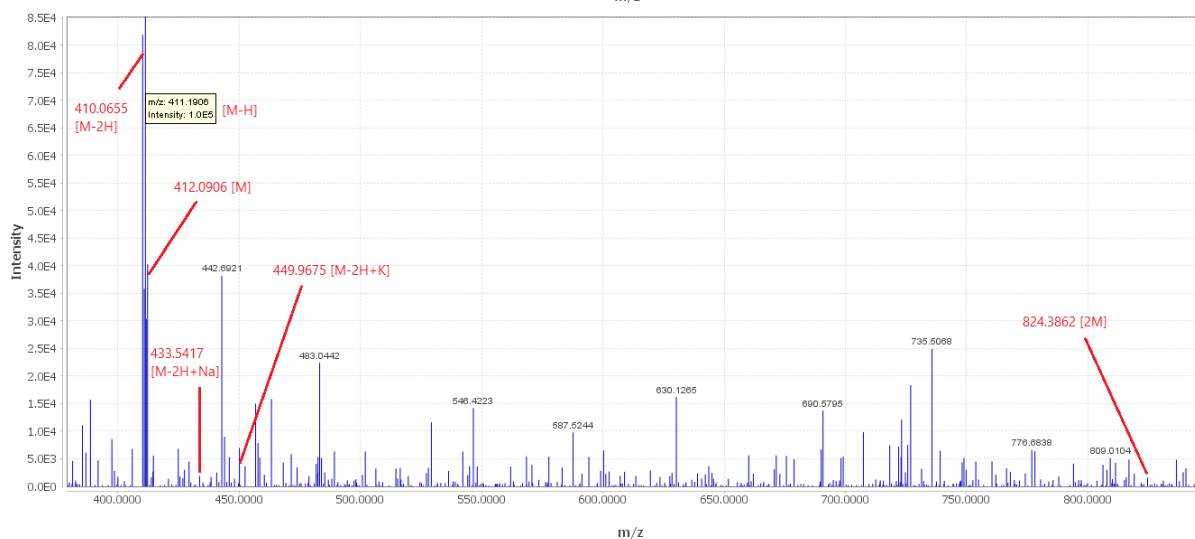
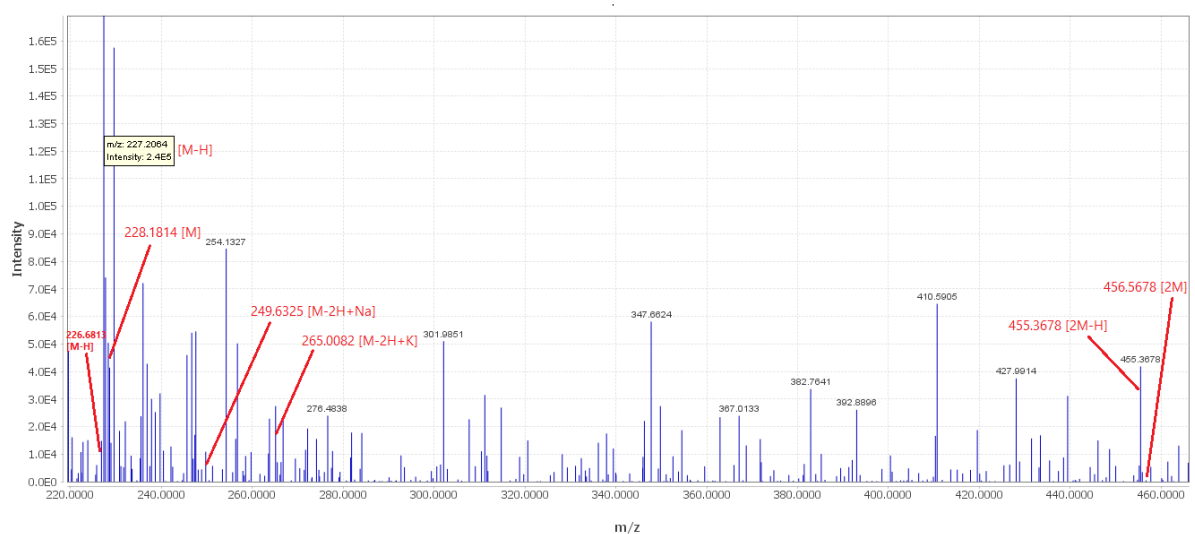
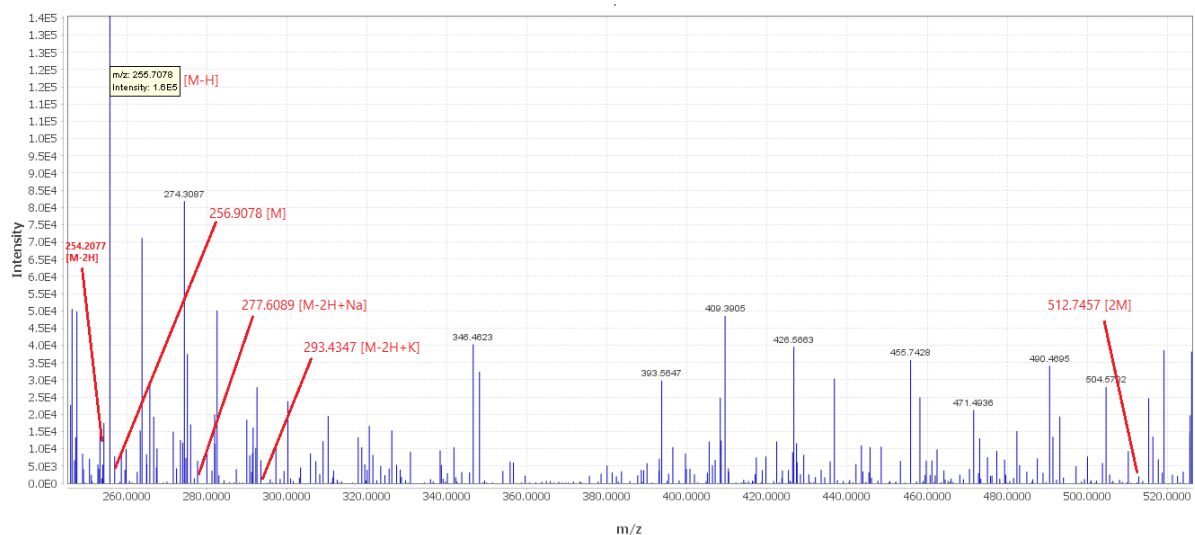


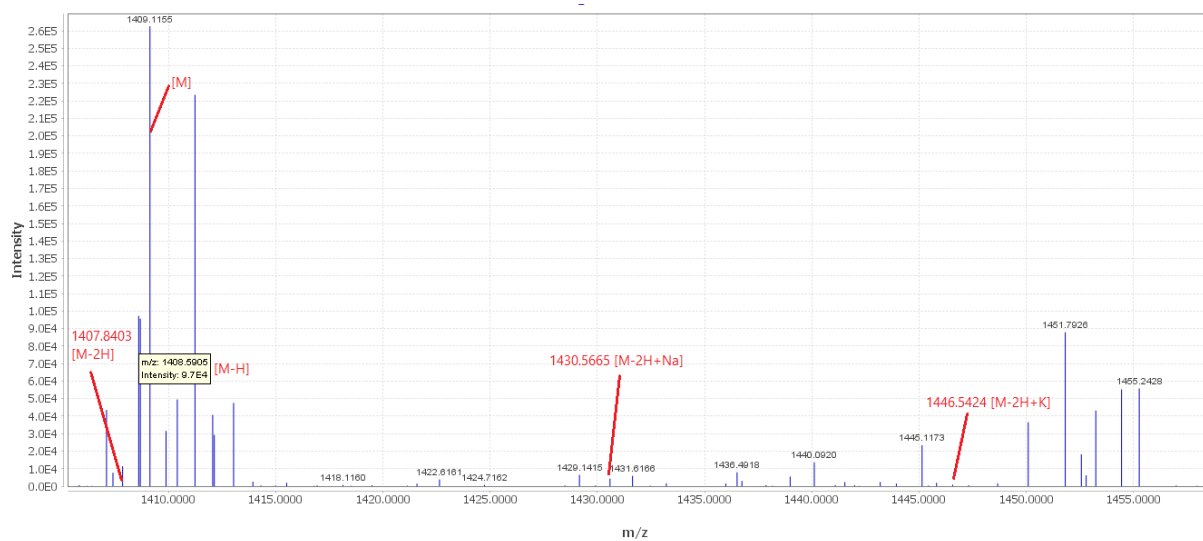
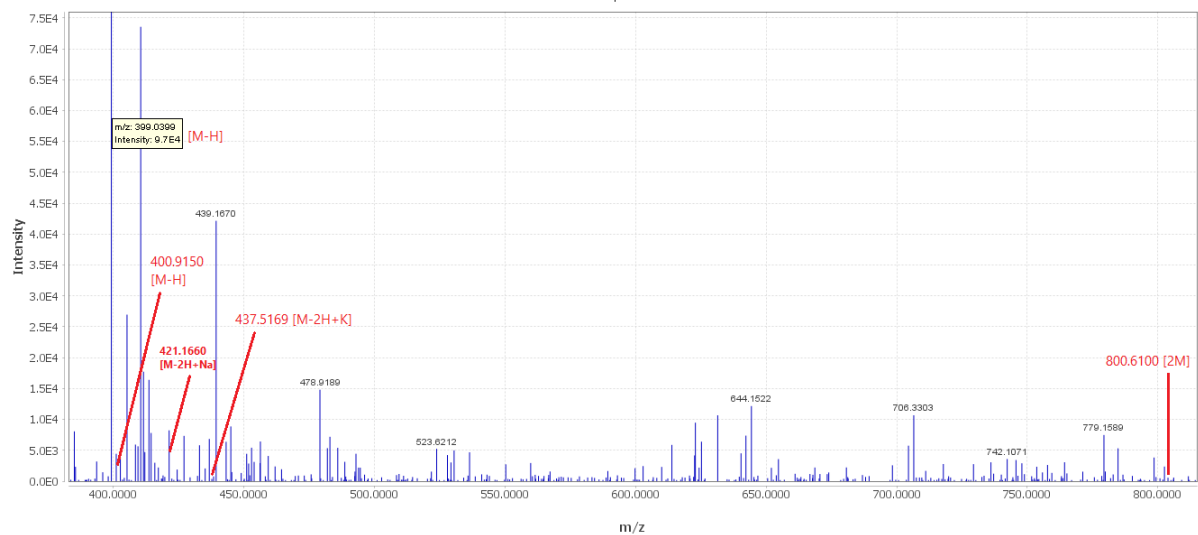


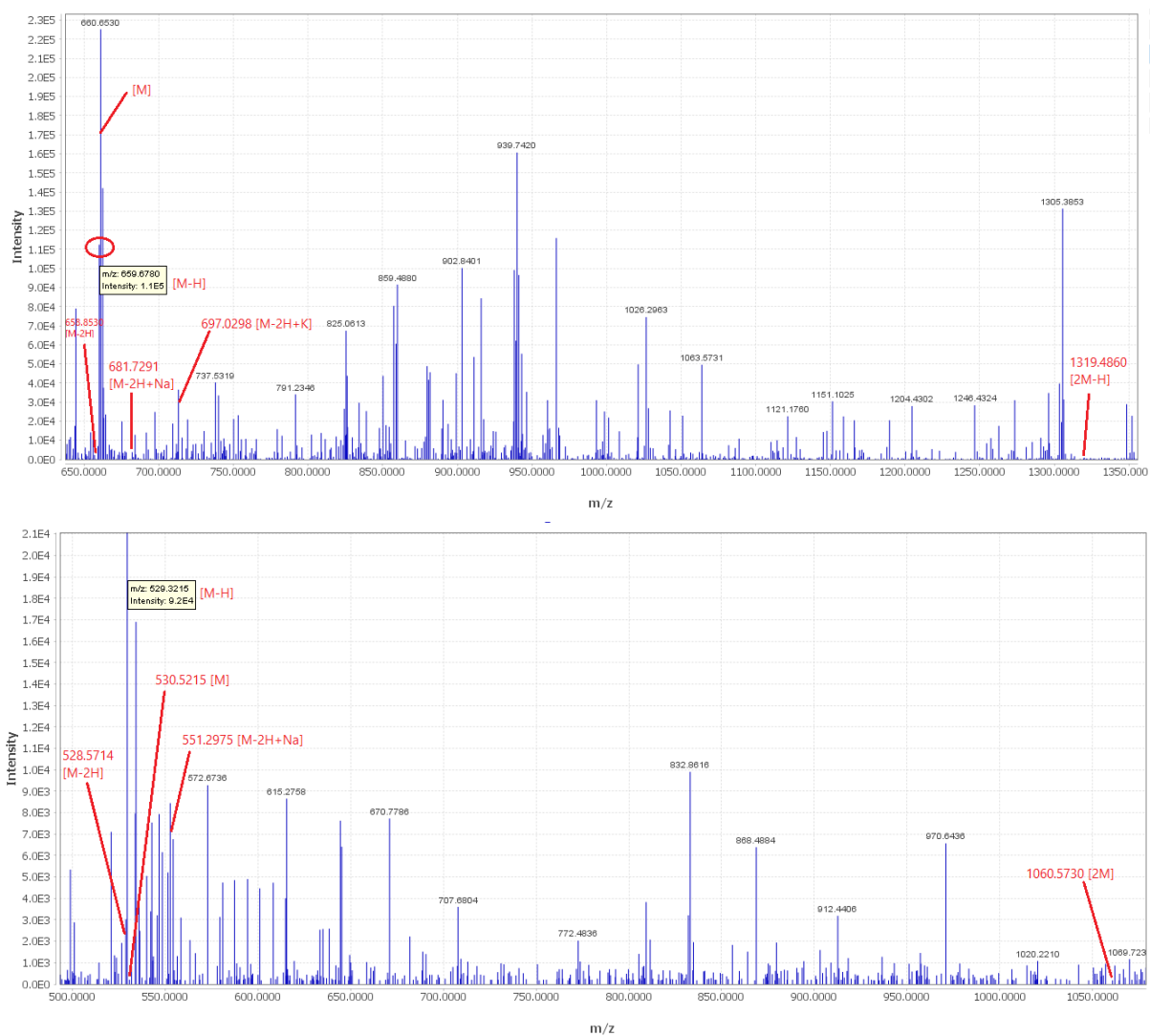












شکل S1- طیف جرمی و اداکت‌های جرمی در ترکیبات شناسایی شده عصاره متانولی ریشه سیلن هرز

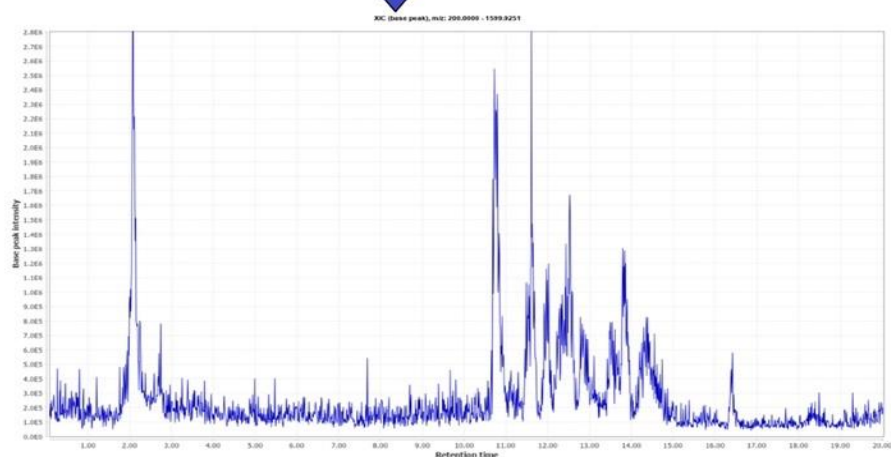
Figure S1- Mass spectrum and mass adducts of identified compounds in methanolic extract of *Silene conoidea* root

(A)

			m/z:		RT:				Identity:		Comment:
ID	Average		Identity	Comment	Peak shape	Minot Galtch 01-08-04 Sample 2.mzML					
ID	m/z	RT				Status	Height	Area			
2491	289.0095	8.59					6.054	4.708			
2492	289.0844	19.26					4.454	6.568			
2493	289.1595	0.06					3.354	6.368			
2494	289.2345	15.14					9.454	7.458			
2495	289.3095	2.14					5.154	6.708			
2496	289.3845	15.94					7.154				
2497	289.4595	3.03					4.254		Show	XIC (base peak) (quick)	
2498	289.5345	0.01					7.054		Search	XIC (dialog)	
2499	289.6095	2.03					6.754		Export	Mass spectrum	
2500	289.6845	4.55					3.554		Identities	Peak in 2D	
2501	289.7595	7.97					6.854		Plot using Intensity Plot module	Peak in 3D	
2502	289.8345	2.58					9.454		Manually define peak	Most Intense MS/MS	
2503	289.9095	13.23					3.254		Delete selected row(s)	All MS/MS	
2504	289.9845	10.86					5.554	8.308	Add new row	MS/MS minor (select 2 rows)	
2505	290.0595	2.08					6.154	9.408		Spectral DB search results	
2506	290.1345	2.15					1.155	1.057		Isotope pattern	
2507	290.2095	2.23					1.365	1.257		Peak row summary	
2508	290.2845	2.27					6.954	1.367			
2509	290.3595	2.17					1.855	1.367			
2510	290.4345	2.21					2.955	1.257			
2511	290.5095	2.31					2.355	1.457			
2512	290.5845	2.17					2.755	1.257			
2513	290.6595	2.22					2.455	1.357			
2514	290.7345	2.20					2.555	1.257			
2515	290.8095	2.44					3.055	1.257			
2516	290.8845	2.23					3.255	1.357			
2517	290.9595	2.24					1.056	2.257			
2518	291.0345	2.21					9.255	2.257			



(B)



(C)

