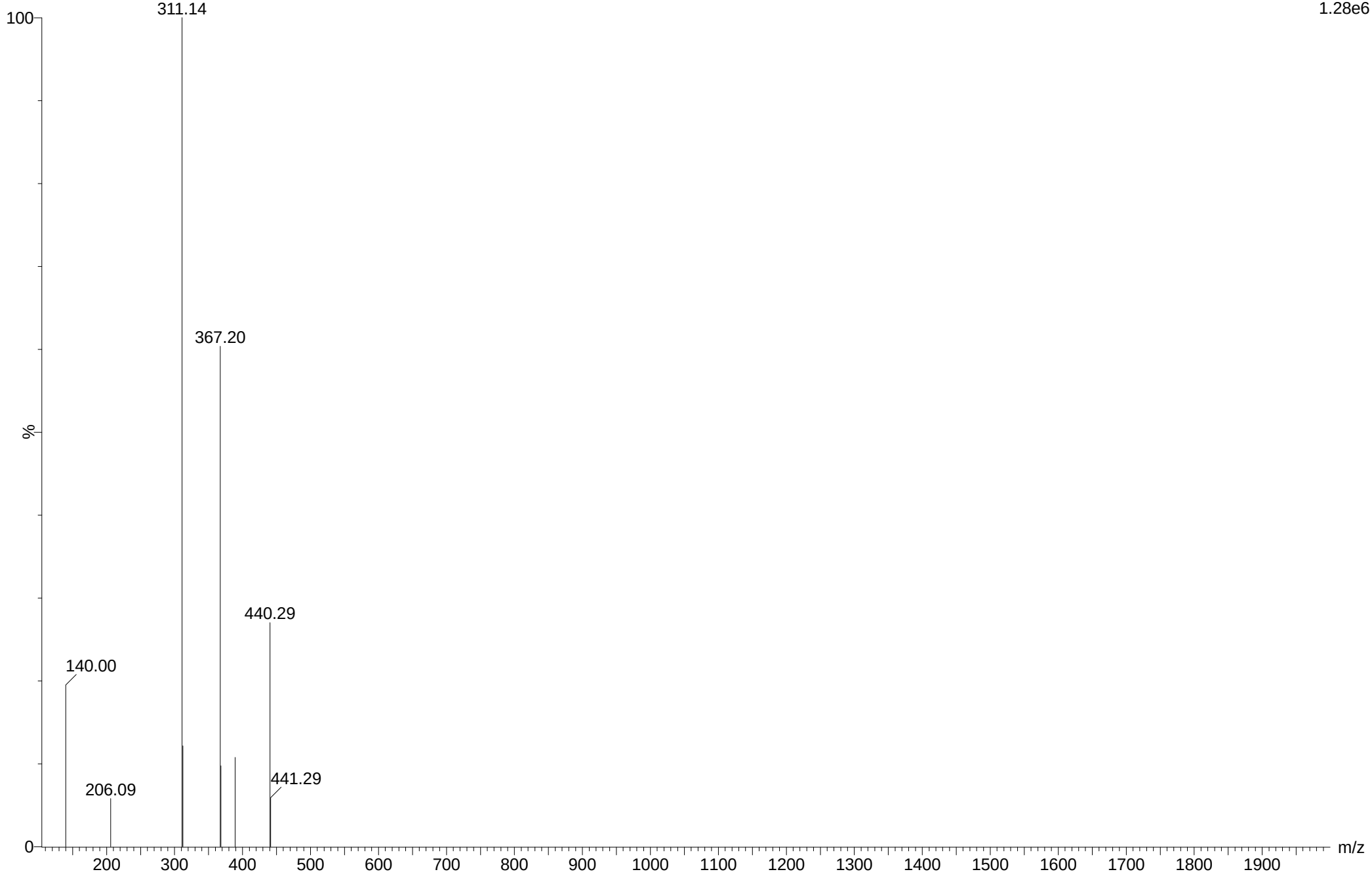
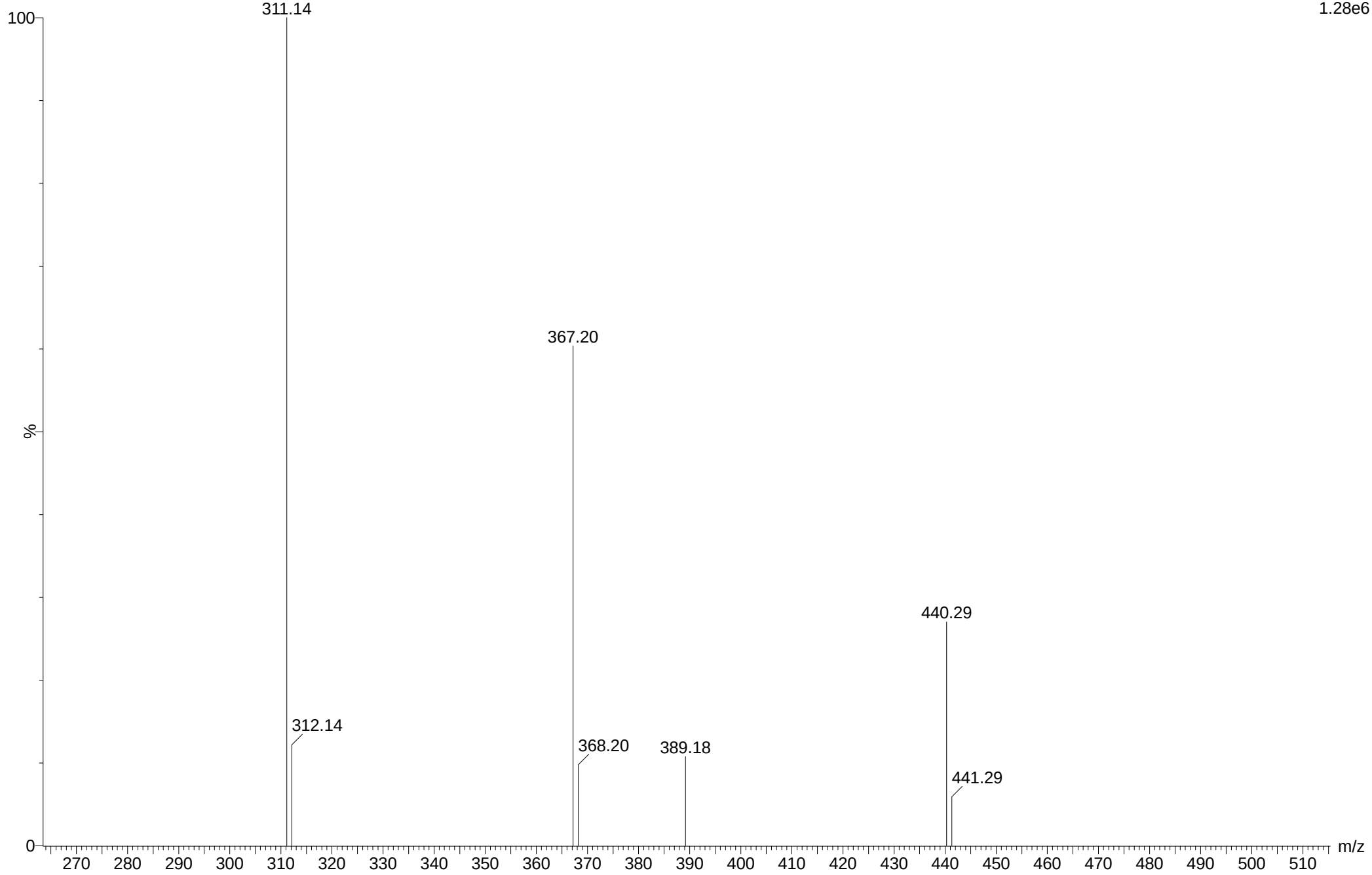
LNB0367-128-P C17H26N4O5 MW=366 University of Birmingham, School of Chemistry Daniel Kovari
Methanol Waters Xevo G2-XS 11-Dec-2020 19:13:35
LRC-DXK-3TVTH-ESI-Pos-1 14 (1.018) AM2 (Ar,30000.0,0.00,0.00); ABS; Cm (14:18) 1: TOF MS ES+
1.28e6



LNB0367-128-P C17H26N4O5 MW=366
Methanol
LRC-DXK-3TVTH-ESI-Pos-1 14 (1.018) AM2 (Ar,30000.0,0.00,0.00); ABS; Cm (14:18)

University of Birmingham, School of Chemistry
Waters Xevo G2-XS

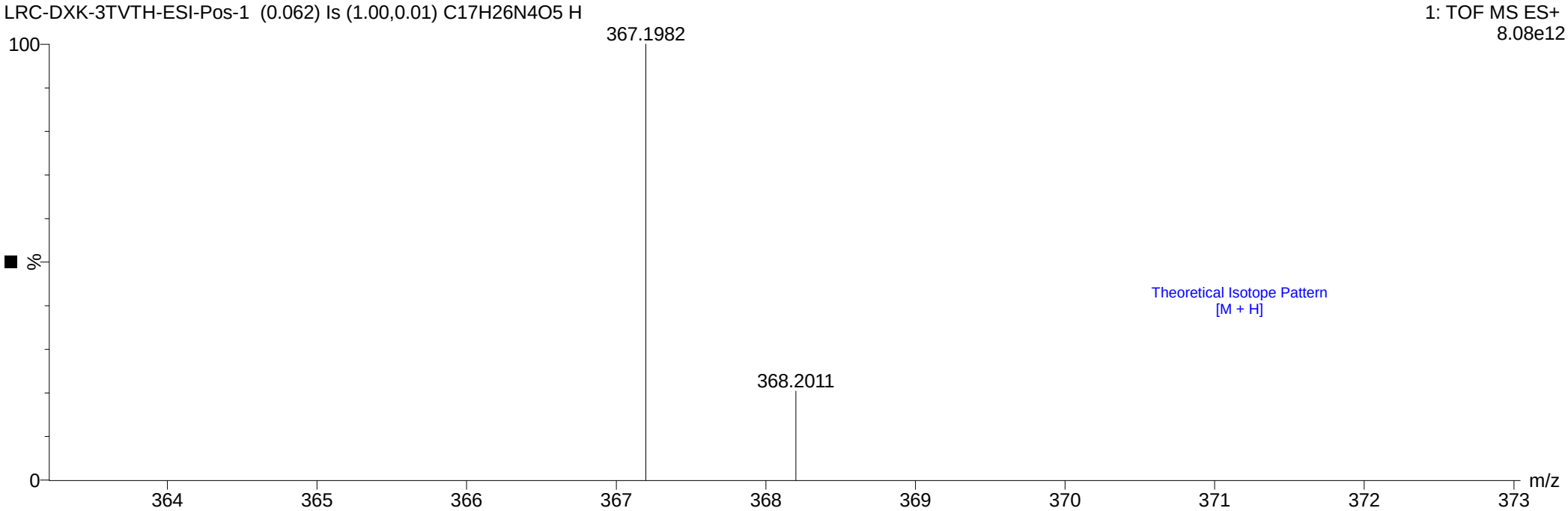
Daniel Kovari
11-Dec-2020 19:13:35
1: TOF MS ES+
1.28e6

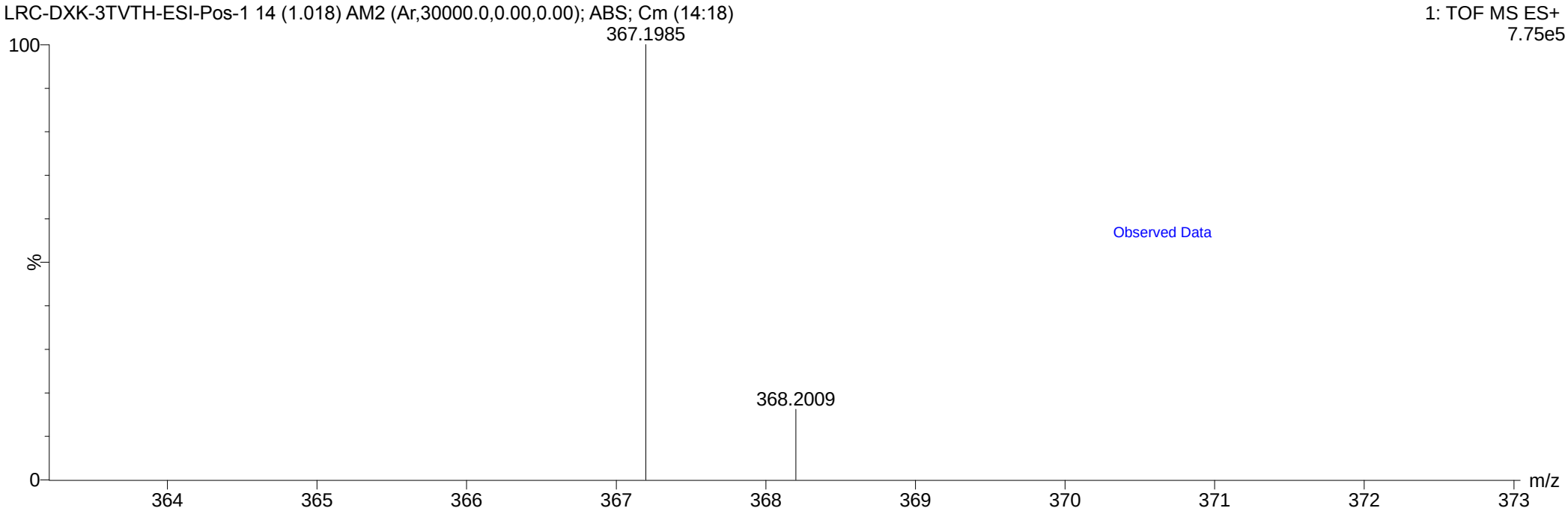


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Single Mass Analysis

Tolerance = 3.0 PPM / DBE: min = -10.0, max = 100.0
Element prediction: Off
Number of isotope peaks used for i-FIT = 4

Monoisotopic Mass, Even Electron Ions
599 formula(e) evaluated with 2 results within limits (all results (up to 1000) for each mass)
Elements Used:
C: 0-50 H: 0-60 N: 0-10 O: 0-10

Minimum:				-10.0				
Maximum:		5.0	3.0	100.0				
Mass	Calc. Mass	mDa	PPM	DBE	i-FIT	Norm	Conf(%)	Formula
367.1985	367.1981	0.4	1.1	6.5	45.9	0.208	81.20	C17 H27 N4 O5
	367.1995	-1.0	-2.7	11.5	47.4	1.671	18.80	C18 H23 N8 O

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Daniel Kovari
11-Dec-2020 19:13:35

LRC-DXK-3TVTH-ESI-Pos-1 14 (1.018) AM2 (Ar,30000.0,0.00,0.00); ABS; Cm (14:18)

1: TOF MS ES+
3.47e5

