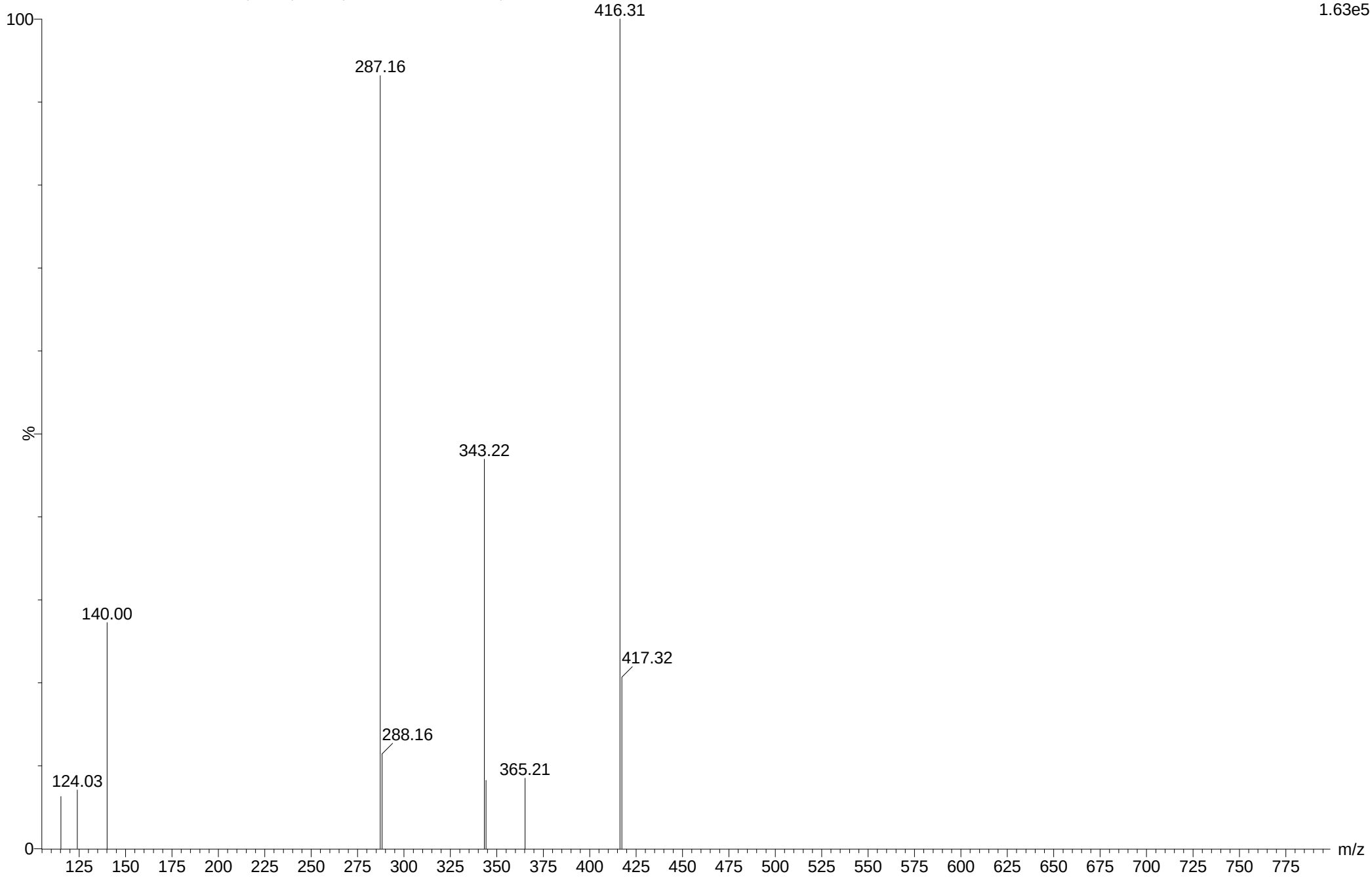
LNB0367-119-P C17H30N2O5 MW=342
Methanol
LRC-DXK-3U44H-ESI-Pos-1 13 (0.922) AM2 (Ar,30000.0,0.00,0.00); ABS

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Waters Xevo G2-XS

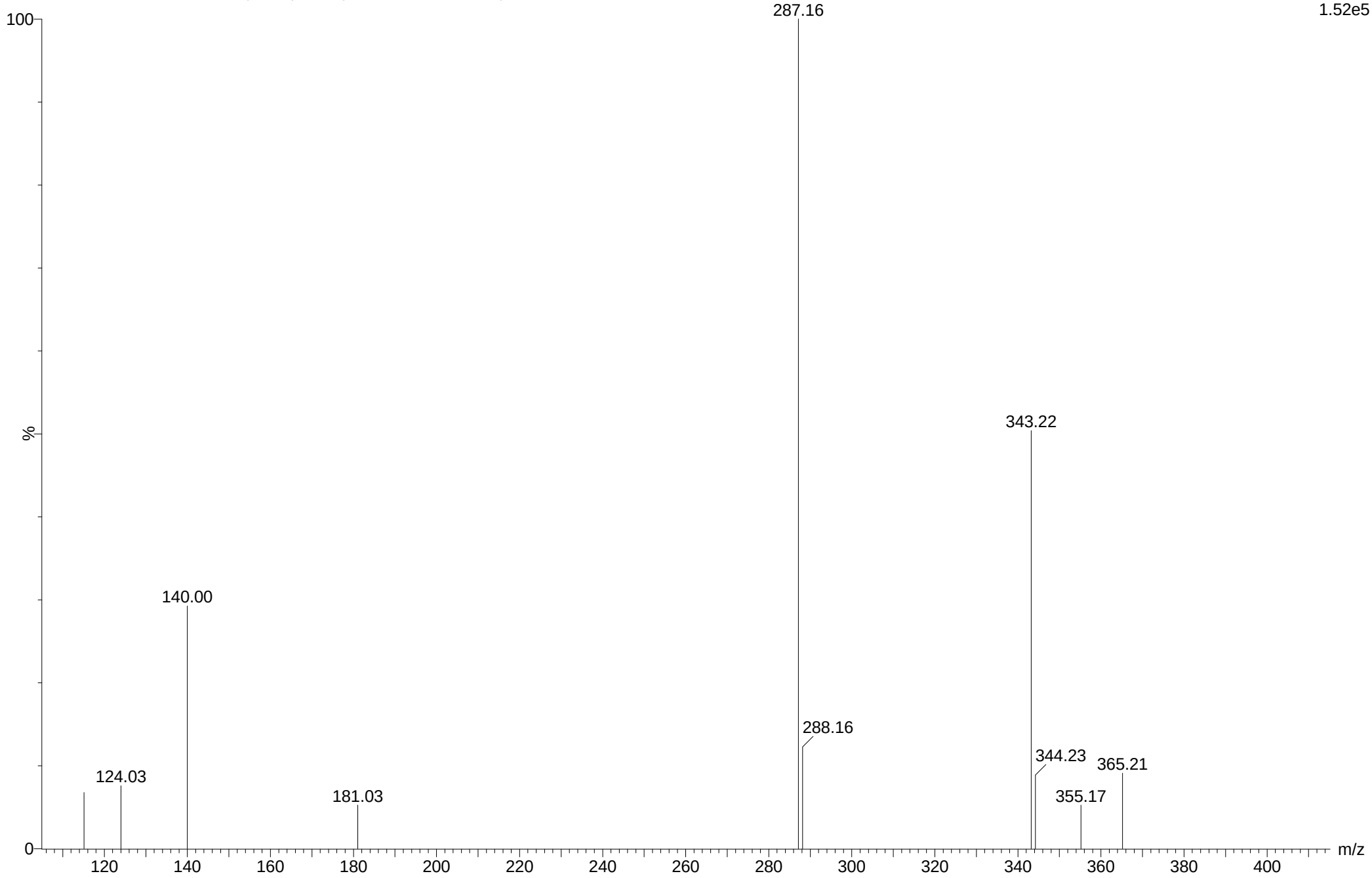
Daniel Kovari
11-Dec-2020 19:31:57
1: TOF MS ES+
1.63e5



LNB0367-119-P C17H30N2O5 MW=342
Methanol
LRC-DXK-3U44H-ESI-Pos-1 13 (0.922) AM2 (Ar,30000.0,0.00,0.00); ABS

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Waters Xevo G2-XS

Daniel Kovari
11-Dec-2020 19:31:57
1: TOF MS ES+
1.52e5



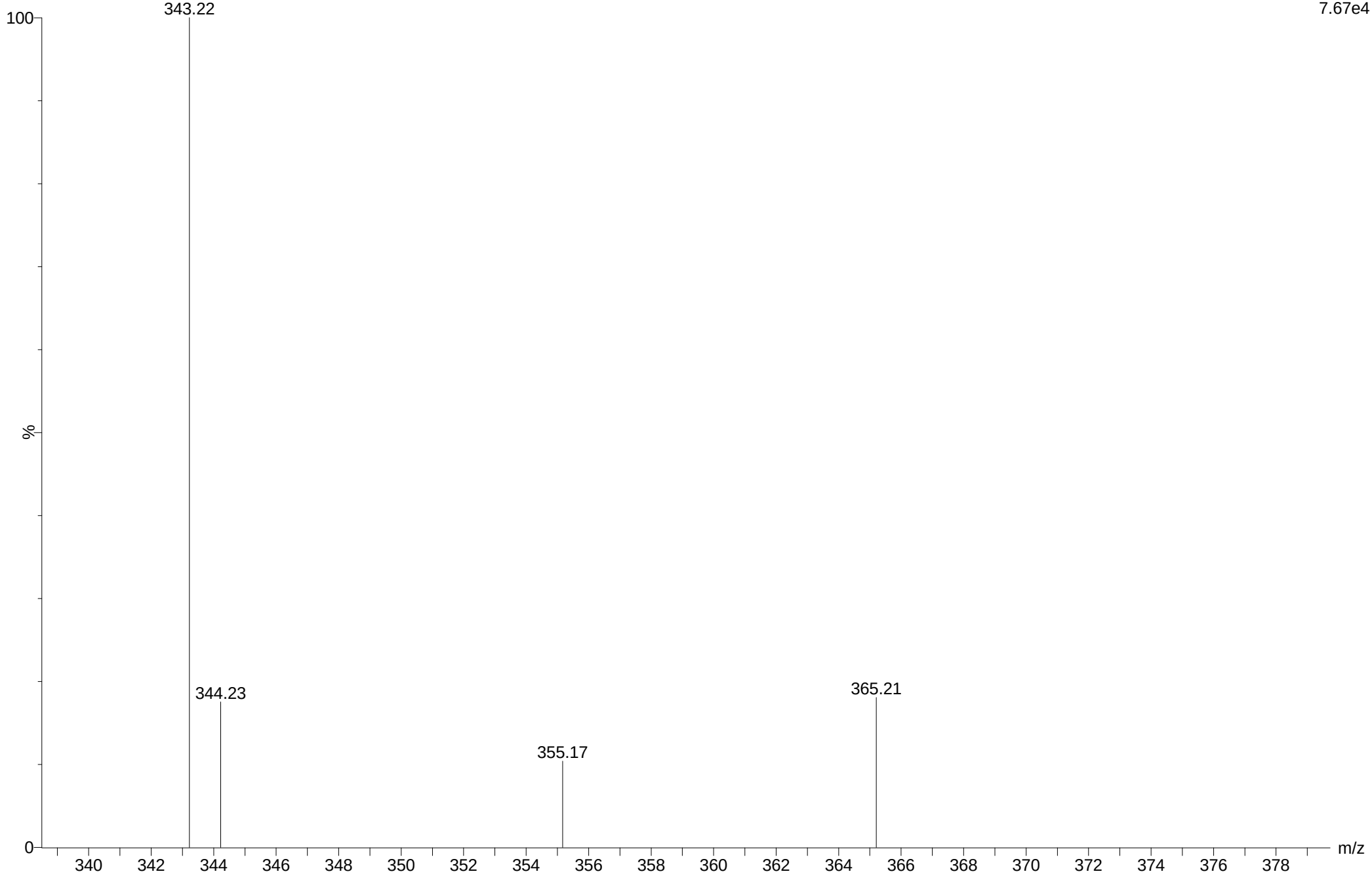
LNB0367-119-P C17H30N2O5 MW=342
Methanol

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

Daniel Kovari
11-Dec-2020 19:31:57

LRC-DXK-3U44H-ESI-Pos-1 13 (0.922) AM2 (Ar,30000.0,0.00,0.00); ABS

1: TOF MS ES+
7.67e4



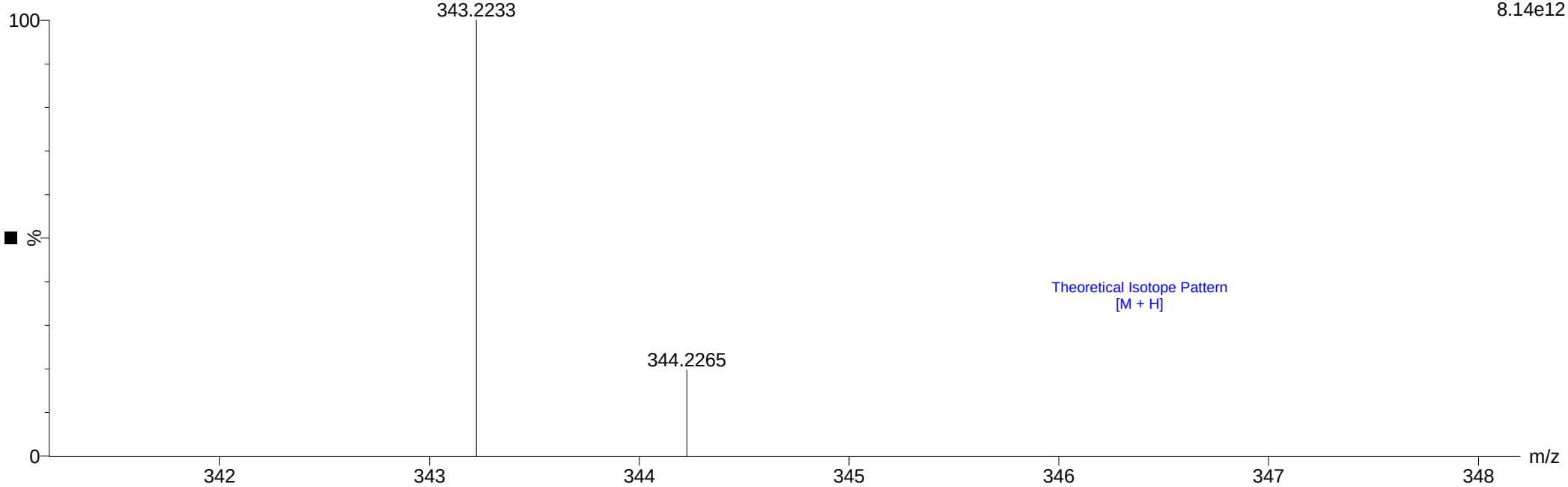
LNB0367-119-P C17H30N2O5 MW=342
Methanol

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

Daniel Kovari
11-Dec-2020 19:31:57

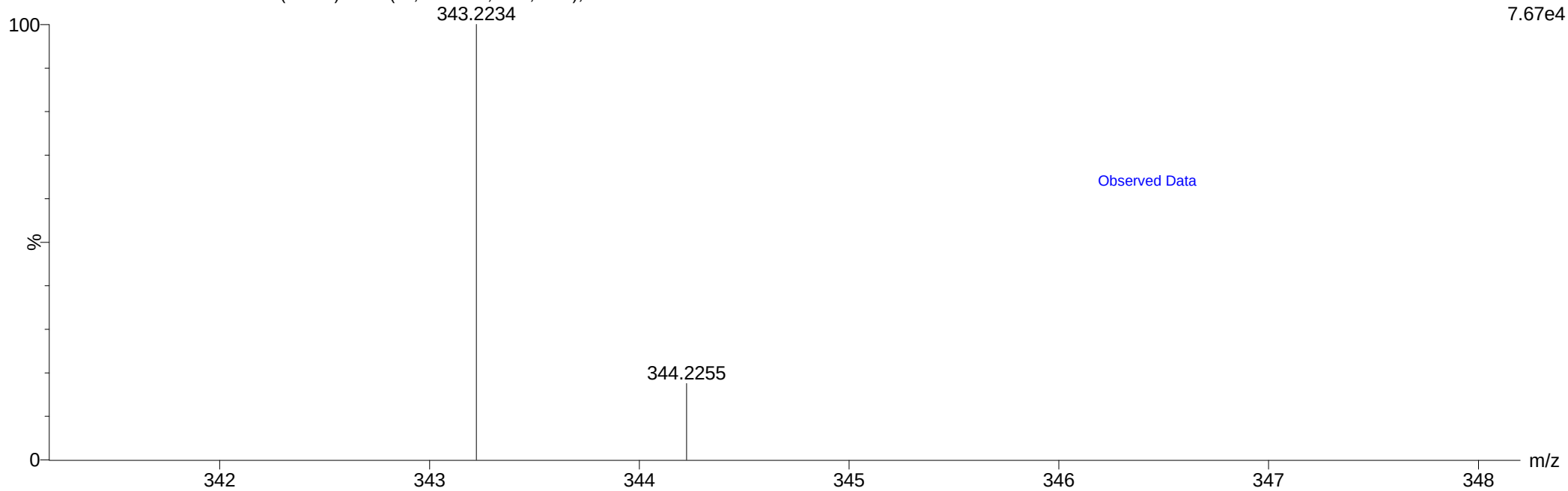
LRC-DXK-3U44H-ESI-Pos-1 (0.062) Is (1.00,0.01) C17H30N2O5 H

1: TOF MS ES+
8.14e12



LRC-DXK-3U44H-ESI-Pos-1 13 (0.922) AM2 (Ar,30000.0,0.00,0.00); ABS

1: TOF MS ES+
7.67e4



Single Mass Analysis

Tolerance = 4.0 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
1629 formula(e) evaluated with 6 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 F: 0-2

Minimum:				-10.0				
Maximum:		5.0	4.0	100.0				
Mass	Calc. Mass	mDa	PPM	DBE	i-FIT	Norm	Conf(%)	Formula
343.2234	343.2233	0.1	0.3	3.5	39.4	2.245	10.59	C17 H31 N2 O5
	343.2237	-0.3	-0.9	8.5	40.3	3.143	4.32	C23 H29 F2
	343.2229	0.5	1.5	-3.5	37.4	0.279	75.67	C7 H29 N8 O5 F2
	343.2225	0.9	2.6	-8.5	42.4	5.316	0.49	C H31 N10 O10
	343.2244	-1.0	-2.9	-0.5	40.5	3.346	3.52	C14 H32 N2 O6 F
	343.2246	-1.2	-3.5	8.5	40.0	2.916	5.41	C18 H27 N6 O