

# Supplementary Materials: Combined (d)SPE-QuEChERS Extraction of Mycotoxins in Mixed Feed Rations and Analysis by High Performance Liquid Chromatography-High-Resolution Mass Spectrometry

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**Table S1.** Regression equations for matrix-matched calibrations of the target mycotoxins in positive mode, using different mobile phase compositions.

Compounds	Ammonium formate 3mM	Ammonium formate 3mM + formic acid 0.1%	Ammonium acetate 3mM	Ammonium acetate 3mM + acetic acid 0.1%
3+15-ADON	$y = 686.6x + 13570$	$y = 632.5x + 20095$	$y = 741.3x - 6619.8$	$y = 575.3x + 23525$
DON	$y = 430.2x - 4372.4$	$y = 364.6x - 5441.5$	$y = 332.9x + 2719.6$	$y = 434.5x - 13436$
ENN B	$y = 2462.5x + 18338$	$y = 783.0x + 27600$	$y = 1897.4x + 14917$	$y = 795.5x + 7306.4$
ENN B <sub>1</sub>	$y = 8491x + 27371$	$y = 3469.3x + 17885$	$y = 6296.3x + 4815.1$	$y = 3413.2x + 24601$
FB <sub>1</sub>	$y = 1819.1x + 15571$	$y = 1629x + 15294$	$y = 1585.2x + 17131$	$y = 1241.4x + 18913$
FB <sub>2</sub>	$y = 1748.4x + 9814$	$y = 1287.5x + 9663.2$	$y = 1770.2x + 3537.9$	$y = 1292.5x + 8063.2$
HT-2	$y = 654.3x - 5538$	$y = 616.1x + 11592$	$y = 546.6x + 11334$	$y = 651.5x + 12973$
MAC A	$y = 5046.2x + 27423$	$y = 3313.5x + 15555$	$y = 4233x + 34351$	$y = 3370.6x + 10198$
OTA	$y = 949.4x + 2287.3$	$y = 1080.5x - 313.4$	$y = 914.4x - 716.28$	$y = 788.2x - 2115.5$
ROQ-C	$y = 2481.1x - 1376.4$	$y = 267.7x + 1894.4$	$y = 2063.3x + 3143.5$	$y = 825.9x - 692.23$
STE	$y = 2474.5x - 49.734$	$y = 2327.9x - 990.45$	$y = 2368.8x - 640.3$	$y = 1988x + 1019.4$
T-2	$y = 4047.8x + 7194.4$	$y = 4099.9x + 9487.3$	$y = 4514.8x + 2602$	$y = 3734.1x - 279.64$
CPA	$y = 1143x - 13643$	$y = 581.1x - 19926$	$y = 1395.3x + 1617.6$	$y = 316.7x - 1866.3$
AND A	$y = 1809.1x + 8567.3$	$y = 970.2x + 9503.1$	$y = 1397x + 6185.6$	$y = 2383.2x - 2540.3$
AOH	$y = 132.2x - 1451.3$	$y = 79.7x + 1439.5$	$y = 124.4x - 1007.5$	$y = 71.3x - 182.9$
MPA	$y = 2150.4x + 945.9$	$y = 4329.2 + 10702$	$y = 2168.3x - 399.3$	$y = 3989.1x + 9532.2$

<b>PEN A</b>	$y = 1096.6x - 672.95$	$y = 671.4x - 1002.7$	$y = 1137.6x + 1878.5$	$y = 702.5x + 1802.5$
<b><math>\alpha</math>-ZOL</b>	$y = 4635.3x + 65293$	$y = 6545.7x + 174983$	$y = 9769x + 21991$	$y = 4653.6x + 148058$
<b><math>\beta</math>-ZOL</b>	$y = 3823x + 52497$	$y = 2121.8x + 63093$	$y = 3326.5x + 20531$	$y = 2151.9x + 63908$
<b>ZEA</b>	$y = 28533x + 23121$	$y = 24249x + 13997$	$y = 30168x + 1236.6$	$y = 23553x + 25766$
<b>FUS X</b>	$y = 318.2x + 17780$	$y = 325.3x + 6505.9$	$y = 476.0x + 2772.7$	$y = 245.8x + 3309.3$
<b>AF B<sub>1</sub></b>	$y = 7189.9x + 3491.5$	$y = 5841.8x + 109544$	$y = 4188.4x + 98650$	$y = 5364.7x + 983.2$
<b>AF B<sub>2</sub></b>	$y = 7472.3x - 5431$	$y = 4863.8x + 31476$	$y = 6580.1x + 3935.8$	$y = 4994.4x - 7080.3$
<b>AF G<sub>1</sub></b>	$y = 5555.5x + 3427.9$	$y = 5117.5x + 5375.1$	$y = 4811.3x + 1885$	$y = 3090x + 5023.2$
<b>AF G<sub>2</sub></b>	$y = 5981x + 42106$	$y = 4215.6x + 33050$	$y = 5137.3x + 15877$	$y = 2822.9x + 23334$

**Table S2.** Regression equations for matrix-matched calibrations of some target mycotoxins in negative mode, using different mobile phase compositions.

Compounds*	Ammonium formate 3mM	Ammonium formate 3mM + formic acid 0.1%	Ammonium acetate 3mM	Ammonium acetate 3mM + acetic acid 0.1%
<b>3+15-ADON</b>	$y = 30.26x + 8404.1$	$y = 13.84x + 349.7$	$y = 331.9x + 22108$	$y = 132.3x + 6379.2$
<b>DON</b>			$y = 106.4x - 2629.5$	$y = 67.92x + 701.6$
<b>FB<sub>1</sub></b>	$y = 2727.1x + 28409$	$y = 1318.5x + 12483$	$y = 2322.4x + 17661$	$y = 1525.9x + 9609.3$
<b>FB<sub>2</sub></b>	$y = 2635.8x + 12000$	$y = 1490.1x + 6757$	$y = 2777.2x + 8039.4$	$y = 1290.7x + 10677$
<b>MAC A</b>			$y = 615.2x + 2850.2$	$y = 381.7x + 36369$
<b>OTA</b>	$y = 2500.6x - 542.8$	$y = 1989.4x - 479.03$	$y = 3107.4x - 9301$	$y = 2487.8x + 419.52$
<b>ROQ-C</b>	$y = 969.9x + 5386.6$	$y = 793.4x + 2059.9$	$y = 1276.1x + 13461$	$y = 585.6x + 5043.5$
<b>STE</b>		$y = 315.3x + 3112.7$		
<b>CPA</b>	$y = 1353x - 86.73$	$y = 817.2x + 8916.4$	$y = 1943.1x + 73304$	$y = 1075.7x + 32130$
<b>AND A</b>	$y = 23861x + 35680$	$y = 27889x + 10102$	$y = 20736x + 6471.2$	$y = 17590x + 17678$
<b>AOH</b>	$y = 1311.8x + 5029.6$	$y = 714.11x + 2167.1$	$y = 1590.4x + 8158.5$	$y = 924.8x + 7135.2$
<b>MPA</b>	$y = 2322.4x + 1374.3$	$y = 2861.6x + 1540.6$	$y = 3856x + 146.64$	$y = 1718.3x + 22040$
<b>PEN A</b>	$y = 1446x + 1127$	$y = 606.9x + 1153.7$	$y = 1576.6x + 1038.4$	$y = 700.77x + 1424.5$
<b>α-ZOL</b>	$y = 44948x + 313726$	$y = 57428x + 169498$	$y = 64727x + 584143$	$y = 38802x + 352390$
<b>β-ZOL</b>	$y = 35230x + 282359$	$y = 17684x + 184993$	$y = 32507x + 230167$	$y = 25063x + 279511$
<b>ZEA</b>	$y = 111740x + 74995$	$y = 83703x + 49378$	$y = 127924x + 79337$	$y = 87939x + 81232$

\*ENN B, ENN B1, HT-2, T-2, FUS X and aflatoxins did not give exploitable calibration curves in negative mode.