

Method Validation For Determination Of Pesticides In Honey

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Abstract—This paper represents the use of ethyl acetate for the determination of pesticides in honey samples by LC-MS/MS detection. The method was validated for 63 pesticides in three different levels. For method validation blank honey samples were spiked at three different levels (10, 20 and 50 µg/kg). The validation parameter included linearity, recovery, accuracy and evaluation of matrix effect. All compounds showed to have good linearity from 5 to 200 µg/kg with a R² of 0.99, in both solvent and matrix calibration curve. The matrix effect did not pass the 20% in none of them. Most of the compound passed the SANTE criteria of the recovery between 70-120 % and RSD lower than 20%.

Keywords—pesticides, honey, SweET, LC-MS/MS

I. INTRODUCTION

The use of pesticides has increased significantly during worldwide. A part the fact that these compounds are beneficiary for the agriculture, many of them reach other then they intended destination. If they are not used properly they can contaminate soil, water and food [1]. In Albania as well as worldwide the use of pesticides in agriculture products has been increased compare to the past. The data has shown that most of the time the farmer are not well aware of the possible contamination from the pesticides uses for their own safety and health [2] as well as for their livestock.

Honey is the substance made of nectar and sweet deposits from plants, gathered, modified and stored in honeycombs by honeybees and as a result it has the image of being natural, healthy and clean. However, today honey is produced in an environment exposed to pollution by different sources of contamination. More specifically, honey can involuntarily be contaminated with pesticides from the environment and also from beekeeping practice [3].

In this work, we showed the validation of a simple method using ethyl acetate [4], which is relatively fast and cost effective, for 63 different pesticides in honey using LC-MS/MS detection system.

II. MATERIAL AND METHODS

Sampling and Sample Preparation

Sample of honey has taken from honey bee farm and well homogenized prior analysis. 10 g of the homogenized honey was accurately weighed in 50 ml polypropylene tubes. The weight samples were diluted with 10 milliliter of distilled water and well homogenized by using a vortex. After homogenization in each samples were added 0.20 mg of PSA (primary secondary amine) and C18 sorbent were added. For the pesticide extraction was used 20 ml of ethyl acetate. The mixture was homogenized by 20 minutes shaking as the first extraction step. A second extraction step was performed for another 10 minutes shaking after 10 g of sodium sulfate was added into the sample. The organic phase was separated by centrifugation and then filter. 1 ml of the filtered organic extract was transferred in 2 ml autosampler vial followed by injection into LC-MS/MS.

2.2. Method Set Up

Pesticides analysis was performed by LC-MS/MS analytical system of Agilent 1200 HPLC system (Agilent Technologies, Germany) with an automatic degasser, a binary pump and an auto sampler connected to the Agilent 6460B Triple-Quad LC/MS system with electrospray ionization interface set at positive and negative polarities. Thermo BDS HYPERSIL C18 Dim.(mm) 150 x 2.1 particle size 3µ. The mobile phase used were: 10-mM ammonium formate (A) and methanol (B) following the gradient program which is shown in table I. The volume of injection was 5µL.

TABLE I. GRADIENT PROGRAM

| | Time | A | B | Flow |
|---|-----------|--------|--------|--------------|
| 1 | 1.00 min | 95.0 % | 5.0 % | 0.550 mL/min |
| 2 | 20.00 min | 5.0 % | 95.0 % | 0.550 mL/min |
| 3 | 27.00 min | 5.0 % | 95.0 % | 0.550 mL/min |
| 4 | 27.20 min | 95.0 % | 5.0 % | 0.550 mL/min |
| 5 | 30.00 min | 95.0 % | 5.0 % | 0.550 mL/min |

Detection of the compounds was performed by Multiple Reaction Monitoring (MRM) with two mass transitions for each pesticide. MRM transitions with their fragmentor voltages (V) and collision energies (CE) are presented in Table II.

TABLE II. RETENTION TIME LC-MS/MS TRANSITIONS, FRAGMENTOR VOLTAGE AND ESI MODE.

| N r. | Cpd Name | RT (min) | Prec Ion | Frag (V) | Quant ion /CE | Qual ion /CE | Mode |
|------|---------------------|----------|----------|----------|---------------|--------------|------|
| 1 | Amitraz | 21.37 | 294.2 | 100 | 163.1 (12) | 122 (30) | ESI+ |
| 2 | Azinophos-Ethyl | 17.71 | 346.1 | 73 | 132.2 (13) | 97.1 (37) | ESI+ |
| 3 | Azinophos-Methyl | 15.5 | 318 | 80 | 160.1 (0) | 132.1 (10) | ESI+ |
| 4 | Azoxystrobine | 16.4 | 404 | 100 | 372 (10) | 344 (23) | ESI+ |
| 5 | Boscaliti | 16.65 | 343.3 | 100 | 307 (15) | 139.9 (20) | ESI+ |
| 6 | Bupiramate | 18.05 | 317.2 | 150 | 166.1 (20) | 108 (25) | ESI+ |
| 7 | Buprofezin | 20.22 | 306 | 100 | 201 (5) | 116.1 (10) | ESI+ |
| 8 | Carbaryl | 13.6 | 202 | 100 | 145 (5) | 127 (25) | ESI+ |
| 9 | Carbendazim | 9.5 | 192.1 | 110 | 160 (18) | 132 (33) | ESI+ |
| 10 | Chlorantraniliprole | 15.73 | 484 | 110 | 453.1(16) | 286.1 (8) | ESI+ |
| 11 | Clofentrezine | 18.95 | 303 | 80 | 138 (10) | 102 (40) | ESI+ |
| 12 | Coumaphos | 18.54 | 363.1 | 140 | 307 (17) | 227 (27) | ESI+ |
| 13 | Cyprodinil | 18.32 | 226 | 135 | 108 (25) | 93 (33) | ESI+ |
| 14 | Difenconazole | 19.44 | 406 | 150 | 251 (25) | - | ESI+ |
| 15 | Diflubenzuron | 17.94 | 309 | 93 | 289.1 (1) | 155.9 (5) | ESI- |
| 16 | Epoxiconazole | 17.69 | 330.1 | 130 | 121.1 (25) | 100.9 (50) | ESI+ |
| 17 | Ethion | 20.29 | 385 | 100 | 199 (5) | 97 (46) | ESI+ |
| 18 | Ethirimol | 14.71 | 210.2 | 120 | 140 (20) | 98 (27) | ESI+ |
| 19 | Ethoprophos | 17.65 | 243.1 | 100 | 173 (10) | 359 (5) | ESI+ |
| 20 | Etofenprox | 22.2 | 394 | 100 | 177 (10) | 10 | ESI+ |
| 21 | Febunconazole | 17.97 | 337 | 150 | 125 (15) | 70 (15) | ESI+ |
| 22 | Fenamidone | 16.55 | 312.2 | 145 | 236.1 (10) | - | ESI+ |
| 23 | Fenarimol | 17.55 | 331 | 150 | 268.1 (20) | 81.1 (30) | ESI+ |
| 24 | Fenpropimorph | 20.4 | 305 | 150 | 147.3 (30) | 117.2 (50) | ESI+ |
| 25 | Fenvalerate | 21.55 | 437 | 90 | 167.1(9) | 125.1 (45) | ESI+ |
| 26 | Fenzaquine | 21.49 | 307 | 120 | 161.1 (15) | 57 (25) | ESI+ |
| 27 | Fipronil | 18.08 | 437 | 130 | 332 (12) | 250 (25) | ESI- |
| 28 | Fipronol desulfinil | 17.79 | 387 | 110 | 351 (10) | 282 (35) | ESI- |
| 29 | Fipronol Sulfon | 18.65 | 453 | 140 | 415 (13) | 282 (25) | ESI- |
| 30 | Fludioxonil | 16.55 | 266 | 90 | 185.1 (20) | 158 (35) | ESI+ |
| 31 | Fluopiram | 17.44 | 397.2 | 83 | 208 (19) | 173 (20) | ESI+ |
| 32 | Fluquiconazole | 17.29 | 376 | 100 | 349.1 (15) | 307.1 (25) | ESI+ |
| 33 | Flusilazole | 18.12 | 316 | 120 | 247 (15) | 165(25) | ESI+ |
| 34 | Hexaconazole | 18.83 | 314 | 130 | 159.1 (40) | 70.1 (20) | ESI+ |
| 35 | Indoxacarb | 19.57 | 528 | 120 | 203 (40) | 150 (20) | ESI+ |
| 36 | Kresoxim-methyl | 18.27 | 314 | 90 | 267 (0) | 116 (10) | ESI+ |
| 37 | Linuron | 15.91 | 248.9 | 110 | 181.9 (10) | 160.1 (15) | ESI+ |
| 38 | Malathion | 16.82 | 331 | 110 | 284 (5) | 127 (5) | ESI+ |
| 39 | Mandipropamid | 16.9 | 412.3 | 110 | 328.2 (8) | 124.9 (30) | ESI+ |
| 40 | Mepanipyrim | 17.21 | 224 | 120 | 106 (25) | 77 (45) | ESI+ |
| 41 | Metaflumizone_A | 19.61 | 505.2 | 150 | 302.1 (15) | 285.1 (45) | ESI- |
| 42 | Metalaxyl | 15.27 | 280.1 | 100 | 220 (10) | - | ESI+ |
| 43 | Methidathion | 15.21 | 303 | 100 | 145 (0) | 85 (15) | ESI+ |
| 44 | Methiocarb | 16.31 | 226 | 100 | 169 (5) | 121 (15) | ESI+ |
| 45 | Methiocarb-Sulphone | 10.29 | 258.1 | 130 | 201.2 (0) | 122.2 (15) | ESI+ |
| 46 | Myclobutanil | 17.18 | 289 | 120 | 125 (40) | 70 (15) | ESI+ |
| 47 | Paclobutrazol | 16.77 | 294.2 | 130 | 125.1 (40) | 70 (15) | ESI+ |
| 48 | Penconazole | 18.4 | 284 | 120 | 159 (30) | 70.1 (15) | ESI+ |
| 49 | Pencycuron | 19.14 | 329.21 | 140 | 218 (5) | 125.0 (25) | ESI+ |
| 50 | Phosmet | 15.56 | 318 | 135 | 160 (22) | 77 (43) | ESI+ |
| 51 | Primicarb | 14.43 | 239.1 | 110 | 182.1 (10) | 72.1 (20) | ESI+ |
| 52 | Prochloraz | 19.05 | 376 | 100 | 308 (5) | 266 (10) | ESI+ |
| 53 | Propiconazol | 18.72 | 342.1 | 140 | 159.2 (25) | 69 (20) | ESI+ |

| | | | | | | | |
|----|----------------|-------|-------|-----|------------|-----------|------|
| 54 | Pyraclostrobin | 18.89 | 388 | 100 | 194.1(5) | 163 (20) | ESI+ |
| 55 | Pyridaben | 20.37 | 322.2 | 100 | 227.1(10) | 96 (10) | ESI+ |
| 56 | Quinoxifen | 20.48 | 308 | 120 | 272 (30) | 197 (35) | ESI+ |
| 57 | Tebuconazole | 18.45 | 308.2 | 140 | 124.9 (35) | 70.1 (20) | ESI+ |
| 58 | Tebuconazole | 18.17 | 353.1 | 80 | 297.1(8) | 133 (20) | ESI+ |
| 59 | Tetraconazole | 17.73 | 372 | 100 | 159.1 (25) | 70.2 (20) | ESI+ |
| 60 | Thiabendazole | 11.12 | 202 | 100 | 175 (24) | 131 (35) | ESI+ |
| 61 | Thiacloprid | 10.61 | 253 | 90 | 126 (20) | 99 (50) | ESI+ |
| 62 | Triadimenol | 16.76 | 296 | 80 | 70 (10) | 43 (35) | ESI+ |
| 63 | Triticonazole | 17.57 | 318.2 | 110 | 125.2 (40) | 70.1 (15) | ESI+ |

III. RESULTS AND DISCUSSION

Calibration Curves and Linearity

Calibration curves of 6 different concentrations were plotted by using matrix matched calibration curve. Concentration range varies from 5-200 µg/kg (using 5,

10, 20, 50, 100 and 200 µg/kg as calibration points). The correlation coefficient (R²) of the calibration curves of all pesticides was ≥0.99. The lowest calibrated level (5 µg/kg) is corresponding the limit of detection (LOD).

TABLE III. CALIBRATION WEIGHT, R² VALUES AND CALIBRATION FORMULAS OF CALIBRATION CURVE IN MATRIX.

| Nr. | Cpd. Name | CF Weight | CF R ² | CF Formula |
|-----|---------------------|-----------|-------------------|-------------------------------------|
| 1 | Amitraz | 1/x | 1.00 | y = 3925.432970 * x - 3479.189550 |
| 2 | Azinophos-Ethyl | 1/x | 1.00 | y = 227.120222 * x - 29.929356 |
| 3 | Azinophos-Methyl | 1/x | 0.99 | y = 921.017769 * x - 1614.907524 |
| 4 | Azoxystrobine | 1/x | 0.99 | y = 11986.041882 * x - 18321.169620 |
| 5 | Boscaliti | 1/x | 0.99 | y = 287.706296 * x - 454.181537 |
| 6 | Bupiramate | 1/x | 0.99 | y = 3207.049127 * x - 5339.291906 |
| 7 | Buprofezin | None | 0.99 | y = 7238.068408 * x - 17390.578834 |
| 8 | Carbaryl | 1/x | 0.99 | y = 3324.118611 * x - 5209.865719 |
| 9 | Carbendazim | None | 1.00 | y = 7698.928313 * x - 9176.023310 |
| 10 | Chlorantraniliprole | 1/x | 0.99 | y = 681.444303 * x - 916.052520 |
| 11 | Clofentrezine | 1/x | 0.99 | y = 267.679288 * x - 381.550935 |
| 12 | Coumaphos | 1/x | 1.00 | y = 1216.399620 * x + 1153.730277 |
| 13 | Cyprodinil | 1/x | 0.99 | y = 1481.906493 * x - 2227.902108 |
| 14 | Difenconazole | 1/x | 0.99 | y = 1001.275982 * x - 1826.995321 |
| 15 | Diflubenzuron | 1/x | 0.99 | y = 416.326481 * x + 16.988936 |
| 16 | Epoxiconazole | 1/x | 0.99 | y = 1428.431848 * x - 2212.036760 |
| 17 | Ethion | None | 0.99 | y = 2136.706635 * x - 4882.152482 |
| 18 | Ethirimol | 1/x | 0.99 | y = 3083.246487 * x - 5226.246593 |
| 19 | Ethoprophos | None | 0.99 | y = 1831.471263 * x - 4007.821222 |
| 20 | Etofenprox | 1/x | 0.99 | y = 1013.861453 * x - 1105.107523 |
| 21 | Febunconazole | 1/x | 0.99 | y = 605.491338 * x - 1100.807523 |
| 22 | Fenamidone | 1/x | 0.99 | y = 1189.302089 * x - 1895.439594 |
| 23 | Fenarimol | None | 0.99 | y = 76.288728 * x - 218.069703 |
| 24 | Fenpropimorph | None | 0.99 | y = 377.555795 * x - 1135.502076 |
| 25 | Fenvalerate | None | 0.99 | y = 7.161244 * x - 9.798213 |
| 26 | Fenzaquine | 1/x | 0.99 | y = 6115.139450 * x - 4696.633730 |
| 27 | Fipronil | 1/x | 0.99 | y = 2001.631008 * x - 2780.456051 |
| 28 | Fipronol desulfinil | None | 1.00 | y = 11020.093444 * x - 8407.639202 |
| 29 | Fipronol Sulfon | None | 1.00 | y = 6159.549343 * x - 3464.896047 |
| 30 | Fludioxonil | 1/x | 0.99 | y = 64.705483 * x - 58.707072 |
| 31 | Fluopiram | 1/x | 0.99 | y = 3422.396525 * x - 5173.598216 |
| 32 | Fluquiconazole | 1/x | 0.99 | y = 82.992716 * x - 57.295303 |
| 33 | Flusilazole | 1/x | 0.99 | y = 933.397852 * x - 1319.947013 |
| 34 | Hexaconazole | 1/x | 0.99 | y = 550.559310 * x - 764.540654 |

| | | | | |
|----|---------------------|------|------|---------------------------------------|
| 35 | Indoxacarb | 1/x | 0.99 | $y = 109.711636 * x - 13.084500$ |
| 36 | Kresoxim-methyl | None | 0.99 | $y = 700.189445 * x - 1592.559793$ |
| 37 | Linuron | None | 0.99 | $y = 29.154845 * x + 62.754474$ |
| 38 | Malathion | 1/x | 0.99 | $y = 2462.687513 * x - 4210.434022$ |
| 39 | Mandipropamid | 1/x | 0.99 | $y = 107.188292 * x - 58.429800$ |
| 40 | Mepanipyrin | 1/x | 0.99 | $y = 1949.208736 * x - 3071.136083$ |
| 41 | Metaflumizone_A | 1/x | 0.99 | $y = 1281.323629 * x - 1928.084453$ |
| 42 | Metalaxyl | 1/x | 0.99 | $y = 338.675214 * x - 91.938048$ |
| 43 | Methidathion | 1/x | 0.99 | $y = 1954.376008 * x - 3091.002452$ |
| 44 | Methiocarb | 1/x | 0.99 | $y = 3566.894243 * x - 5424.704330$ |
| 45 | Methiocarb-Sulphone | 1/x | 0.99 | $y = 168.306655 * x - 232.404110$ |
| 46 | Myclobutanil | 1/x | 0.99 | $y = 775.865216 * x - 1019.725373$ |
| 47 | Paclobutrazol | 1/x | 0.99 | $y = 1656.280328 * x - 2276.221347$ |
| 48 | Penconazole | 1/x | 0.99 | $y = 1324.807288 * x - 1538.257694$ |
| 49 | Pencycuron | 1/x | 0.99 | $y = 339.352933 * x - 305.995487$ |
| 50 | Phosmet | None | 0.99 | $y = 28.271880 * x - 114.973685$ |
| 51 | Primicarb | None | 1.00 | $y = 14712.070106 * x - 54258.981409$ |
| 52 | Prochloraz | 1/x | 0.99 | $y = 107.679339 * x - 68.566912$ |
| 53 | Propiconazol | 1/x | 0.99 | $y = 791.274055 * x - 909.351547$ |
| 54 | Pyraclostrobin | 1/x | 0.99 | $y = 3885.615936 * x - 5635.817691$ |
| 55 | Pyridaben | 1/x | 0.99 | $y = 7467.848775 * x - 11461.127197$ |
| 56 | Quinoxifen | 1/x | 0.99 | $y = 503.217577 * x - 711.571170$ |
| 57 | Tebuconazole | 1/x | 0.99 | $y = 1555.420127 * x - 1812.516124$ |
| 58 | Tebufenozide | 1/x | 0.99 | $y = 281.545533 * x - 421.717974$ |
| 59 | Tetraconazole | 1/x | 0.99 | $y = 309.648251 * x - 401.284151$ |
| 60 | Thiabendazole | 1/x | 0.99 | $y = 203.351918 * x - 263.440544$ |
| 61 | Thiacloprid | 1/x | 0.99 | $y = 162.441346 * x - 67.979064$ |
| 62 | Triadimenol | 1/x | 0.99 | $y = 233.071926 * x - 267.331958$ |
| 63 | Triticonazole | None | 0.99 | $y = 74.779013 * x - 15.099546$ |

Matrix Effects

Matrix effect was Calculated based on slopes of calibration curves following the above formula:

$$\%ME = \frac{\text{Slope Of MMC Curve}}{\text{Slope SMC Curve}} * 100$$

From the percentage of matrix effect is given in table III.

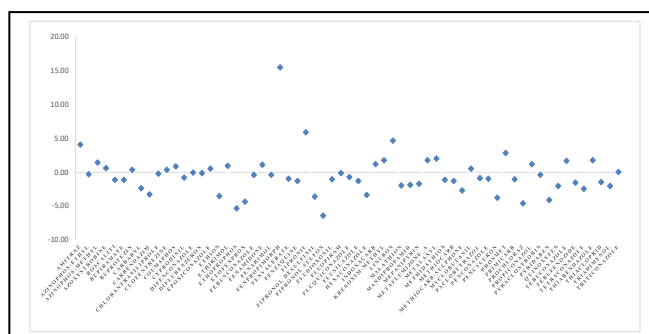


Fig. 1. Matrix effect plot for all the pesticides

None from the 63 compounds represented in this study shows to have matrix effect lower than 20 %, this can be due to the dilution of sample before analysis [5].

Precision in Spiked Matrix Samples

Method performance was evaluated by checking two parameters the accuracy and precision according to SANTE guidance 2017[6].

In order to study the recovery, prior extraction the honey samples were spiked in three different levels (10, 20 and 50 µg/kg). Number of replicates used were six for each spiking level, the recovery and RSD was calculated for each level. The analytical performance parameters for all the pesticides tested in this study using LC-MS/MS system are given in table III.

In the figure 2 are giving the recovery measurements versus RSD in percentage, for each recovery level.

In the first spiking level of 10 µg/kg 59 out of 63 - pesticides have met the recovery criteria between 70-120 % and the RSD <20%. Although 3 compounds (fenvalerate, linuron, triticonazole) have the recovery between 70 and 120 % their RSD was very high, which means that these three compounds did not meet the SANTE criteria for recovery and RSD. On the other hand, Amitraz had a recovery of 34 % but its RSD was 3.9, which according to SANTE guidance is accepted. Ethionphos did not pass this validation level because it has a low recovery as well as a high RSD.

TABLE IV. AVERAGE RECOVERIES (N = 6), AND RSD FOR SAMPLES FORTIFIED AT 10, 20, 50 µG/KG

| Nr. | Cpd. Name | Spiking level | | Spiking level | | Spiking level | | |
|-----|---------------------|---------------|--------|---------------|--------|---------------|--------|-------|
| | | 0.01 | mg/kg | 0.02 | mg/kg | 0.05 | mg/kg | |
| | | Recovery% | RSDr % | Recovery% | RSDr % | Recovery % | RSDr % | |
| 1 | Amitraz | 39.1 | 3.9 | 47.9 | 7.48 | 47.9 | 7.5 | 4.04 |
| 2 | Azinophos-Ethyl | 79.4 | 11.7 | 82.5 | 5.0 | 85.6 | 6.4 | -0.39 |
| 3 | Azinophos-Methyl | 99.2 | 2.7 | 92.2 | 3.0 | 87.8 | 3.3 | 1.38 |
| 4 | Azoxystrobin | 91.4 | 2.9 | 89.8 | 2.9 | 89.6 | 3.7 | 0.52 |
| 5 | Boscaliti | 100.2 | 3.4 | 94.8 | 7.9 | 86.7 | 3.3 | -1.24 |
| 6 | Bupiramate | 86.8 | 3.5 | 85.4 | 3.4 | 81.5 | 5.4 | -1.21 |
| 7 | Buprofezin | 90.7 | 3.7 | 84.6 | 5.5 | 75.8 | 5.8 | 0.27 |
| 8 | Carbaryl | 89.6 | 2.3 | 90.3 | 3.3 | 90.4 | 2.6 | -2.46 |
| 9 | Carbendazim | 91.8 | 4.1 | 90.0 | 0.7 | 88.6 | 1.5 | -3.32 |
| 10 | Chlorantraniliprole | 94.9 | 2.6 | 90.6 | 5.1 | 89.7 | 3.8 | -0.32 |
| 11 | Clofentrezine | 86.3 | 11.8 | 87.0 | 5.6 | 79.1 | 5.7 | 0.32 |
| 12 | Coumaphos | 92.1 | 5.6 | 81.8 | 9.3 | 73.7 | 5.0 | 0.80 |
| 13 | Cyprodinil | 91.4 | 4.4 | 86.9 | 4.7 | 84.6 | 6.1 | -0.91 |
| 14 | Difenconazole | 93.5 | 5.0 | 87.0 | 3.2 | 79.4 | 5.1 | -0.11 |
| 15 | Diflubenzuron | 84.0 | 13.8 | 89.6 | 7.3 | 93.0 | 3.3 | -0.19 |
| 16 | Epoconazole | 100.7 | 2.8 | 95.3 | 4.6 | 88 | 2.93 | 0.46 |
| 17 | Ethion | 86.5 | 2.9 | 83.3 | 7.5 | 107.1 | 6.4 | -3.59 |
| 18 | Ethirimol | 93.6 | 1.3 | 91.9 | 3.0 | 89.2 | 1.8 | 0.85 |
| 19 | Ethoprophos | 39.4 | 21.5 | 87.8 | 7.8 | 115.6 | 7.4 | -5.42 |
| 20 | Etofenprox | 91.1 | 8.7 | 87.8 | 7.8 | 65.9 | 6.1 | -4.43 |
| 21 | Febunconazole | 95.7 | 5.8 | 91.7 | 6.2 | 87.3 | 4.2 | -0.43 |
| 22 | Fenamidone | 93.4 | 3.9 | 88.3 | 6.9 | 87.2 | 4.3 | 1.04 |
| 23 | Fenarimol | 101.7 | 11.5 | 104.5 | 9.8 | 91.7 | 10.4 | -0.48 |
| 24 | Fenpropimorph | 77.3 | 12.8 | 77.8 | 11.1 | 67.5 | 5.6 | 15.50 |
| 25 | Fenvalerate | 76.1 | 43.7 | 77.6 | 28.5 | 84.0 | 17.5 | -0.99 |
| 26 | Fenzaquine | 81.8 | 7.1 | 81.7 | 5.2 | 76.9 | 6.7 | -1.40 |
| 27 | Fipronil | 88.2 | 3.0 | 85.6 | 2.7 | 79.5 | 4.4 | 5.89 |
| 28 | Fipronol desulfinil | 78.3 | 6.4 | 78.2 | 8.4 | 74.1 | 3.8 | -3.66 |
| 29 | Fipronol Sulfon | 73.6 | 7.0 | 75.4 | 7.8 | 70.0 | 5.3 | -6.51 |
| 30 | Fludioxonil | 85.1 | 13.8 | 89.9 | 7.7 | 84.3 | 7.8 | -1.11 |
| 31 | Fluopiram | 90.7 | 2.5 | 87.9 | 2.8 | 85.2 | 3.1 | -0.16 |
| 32 | Fluquinconazole | 88.4 | 19.9 | 88.5 | 8.9 | 89.4 | 6.9 | -0.79 |
| 33 | Flusilazole | 93.4 | 8.4 | 92.6 | 5.9 | 85.2 | 4.6 | -1.39 |
| 34 | Hexaconazole | 94.9 | 10.2 | 90.0 | 3.6 | 87.3 | 6.3 | -3.45 |
| 35 | Indoxacarb | 82.0 | 10.9 | 79.7 | 8.2 | 76.5 | 11.6 | 1.08 |
| 36 | Kresoxim-methyl | 92.1 | 3.8 | 85.6 | 2.3 | 82.9 | 6.2 | 1.69 |
| 37 | Linuron | 105.7 | 52.4 | 69.1 | 10.8 | 82.3 | 19.0 | 4.60 |
| 38 | Malathion | 75.5 | 7.8 | 83.0 | 3.1 | 84.2 | 2.4 | -1.99 |
| 39 | Mandipropamid | 88.9 | 4.9 | 89.2 | 8.7 | 86.0 | 5.8 | -1.92 |
| 40 | Mepanipyrim | 93.0 | 3.4 | 90.3 | 2.3 | 87.1 | 3.1 | -1.81 |
| 41 | Metaflumizone_A | 89.0 | 3.8 | 82.9 | 6.9 | 74.5 | 7.8 | 1.67 |
| 42 | Metalaxyl | 86.8 | 5.5 | 94.9 | 2.0 | 91.6 | 2.6 | 1.96 |
| 43 | Methidathion | 87.5 | 3.1 | 89.6 | 2.5 | 88.4 | 2.7 | -1.17 |
| 44 | Methiocarb | 83.4 | 3.7 | 85.8 | 2.9 | 87.0 | 3.6 | -1.32 |
| 45 | Methiocarb-Sulphone | 77.1 | 4.4 | 82.6 | 3.9 | 82.9 | 4.4 | -2.79 |
| 46 | Myclobutanil | 93.9 | 7.4 | 91.4 | 2.5 | 92.2 | 3.6 | 0.48 |
| 47 | Paclobutrazol | 89.6 | 5.6 | 90.6 | 3.2 | 89.4 | 4.3 | -0.96 |
| 48 | Penconazole | 93.9 | 4.6 | 92.9 | 3.0 | 86.4 | 3.6 | -1.05 |
| 49 | Pencycuron | 92.8 | 11.3 | 87.7 | 12.4 | 77.4 | 4.4 | -3.87 |
| 50 | Phosmet | 109 | 16.04 | 89.4 | 8.5 | 82.1 | 7.3 | 2.76 |
| 51 | Primicarb | 99.3 | 2.4 | 96.2 | 1.5 | 89.4 | 2.2 | -1.13 |
| 52 | Prochloraz | 110.6 | 17.7 | 87.2 | 14.2 | 89.7 | 8.6 | -4.64 |
| 53 | Propiconazol | 94.4 | 5.5 | 90.4 | 7.1 | 88.5 | 4.2 | 1.13 |
| 54 | Pyraclostrobin | 86.8 | 4.6 | 83.0 | 5.5 | 76.8 | 5.9 | -0.44 |
| 55 | Pyridaben | 76.2 | 4.7 | 75.0 | 6.6 | 70.0 | 4.98 | -4.14 |

| | | | | | | | | |
|----|---------------|-------|------|------|------|------|------|-------|
| 56 | Quinoxifen | 81.8 | 8.4 | 79.9 | 4.9 | 72.5 | 7.0 | -2.08 |
| 57 | Tebuconazole | 97.6 | 3.0 | 95.5 | 3.6 | 87.9 | 5.9 | 1.65 |
| 58 | Tebuconazole | 97.1 | 8.8 | 88.3 | 8.3 | 87.5 | 13.6 | -1.62 |
| 59 | Tetraconazole | 91.5 | 10.5 | 91.3 | 8.0 | 83.6 | 7.1 | -2.55 |
| 60 | Thiabendazole | 102.1 | 2.0 | 98.3 | 4.1 | 93.9 | 2.7 | 1.67 |
| 61 | Thiacloprid | 97.9 | 2.5 | 98.4 | 3.5 | 93.0 | 2.8 | -1.52 |
| 62 | Triadimenol | 89.0 | 4.2 | 92.3 | 6.3 | 91.1 | 4.6 | -2.14 |
| 63 | Triticonazole | 82.5 | 23.3 | 92.7 | 15.0 | 94.8 | 13.5 | -0.01 |

The second level of validation was 20 µg/kg. In this level only one compound, fenvalerate and linuron, did not met the SANTE [6], and amitraz still has the same tendency of low recovery and low RSD.

All the compound did pass the criteria of having a recovery between 70-120 % and RSD lower than 20 %. Amitraz still have the same tendency even in this spiking level to have low recovery and low RSD.

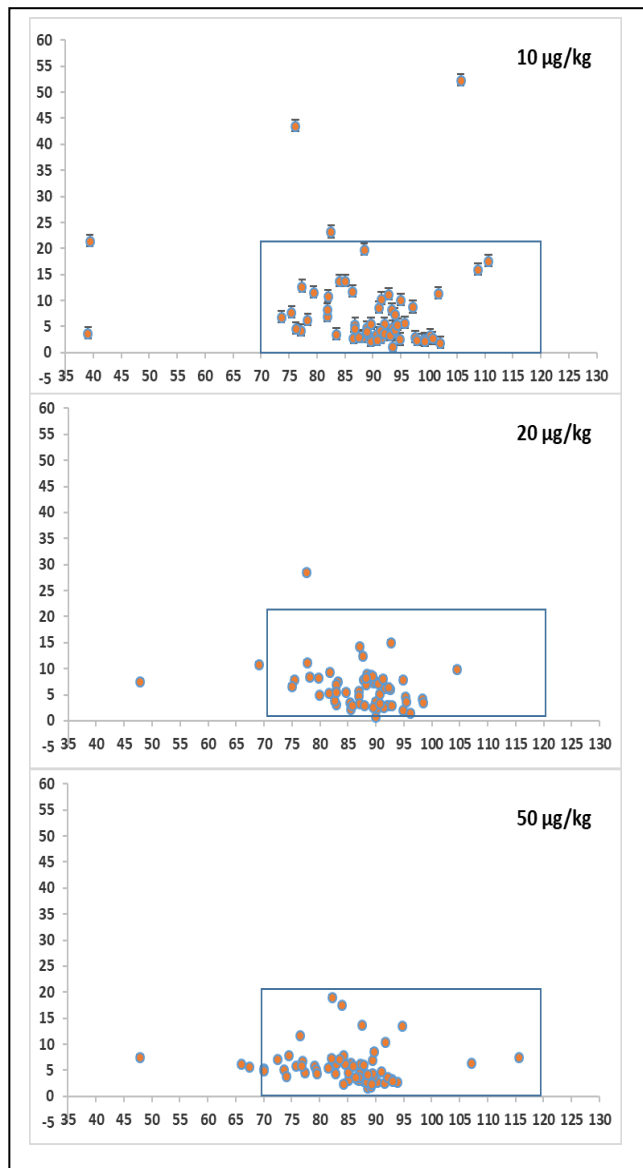


Fig. 2. Example of a figure caption. (figure caption)

IV. CONCLUSION

In this study we presents a reliable and cost effective multi residue method for the determination of different pesticides in honey, which can be either used in

apiculture or in the surrounding agriculture in the context food safety study. The procedure includes honey dilution with distillate water and the ethyl acetate extraction followed by analytical measurement in LC-MS/MS without further clean up procedure. The method that we validated complies with the validation requirements of the SANTE document.

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