

**Supplementary Table 3.** Description of the processes and parameters applied to LC-MS chromatograms with MZmine ver.2.38 to obtain the metabolomic fingerprints of *Arabidopsis* root samples from both positive and negative ionization modes.

		Parameters
<b>1</b>	<b>Baseline correction – RollingBall baseline corrector</b>	
	Chromatogram type	TIC
	Use m/z bins	No
	wm	25
	ws	25
<b>2</b>	<b>Mass detection (exact Mass)</b>	
	Noise level	$1 \times 10^3$
<b>3</b>	<b>FTMS shoulder peak filter</b>	
	Mass Resolution	60,000
	Peak model function	Lorentzian
<b>4</b>	<b>Chromatogram builder</b>	
	Minimum time span	0.04
	Minimum height	$1 \times 10^3$
	m/z tolerance	0.0005 m/z or 7ppm
<b>5</b>	<b>Smoothing</b>	
	Filter width	5
<b>6</b>	<b>Chromatogram deconvolution (local minimum search)</b>	
	Chromatographic threshold	40%
	Search minimum in RT range (min)	0.3
	Minimum relative height	30%
	Minimum absolute height	3000
	Minimum ratio of peak top/edge	1.5
	Peak duration range	0-1
<b>7</b>	<b>Normalization – Retention time normalizer</b>	
	m/z tolerance	0.0005 m/z or 7ppm
	RT tolerance	0.3
	Minimum Standard Intensity	$1 \times 10^5$
<b>8</b>	<b>Chromatogram alignment (join alignment)</b>	
	m/z tolerance	0.0005 m/z or 6ppm
	Weight for m/z	80
	RT tolerance	0.25
	Weight for RT	40
<b>9</b>	<b>Gap filling (Peak Finder)</b>	
	Intensity tolerance	30%
	m/z tolerance	0.0005 m/z or 6ppm
	Retention time tolerance	0.25
	RT correction	Yes
<b>10</b>	<b>Metabolite Assignment</b>	
	m/z tolerance	0.001 m/z or 15ppm*
	RT tolerance	0.45*
<b>11</b>	<b>Data Exported</b>	Peak Area

RT, retention time; m/z, mass to charge ratio

\* Metabolite assignation filtered a posteriori (Supplementary Table 5)