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## Submerged macrophyte assessment in rivers: An automatic mapping method using Pléiades imagery

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20 **Abstract:**

21 Submerged macrophyte monitoring is a major concern for hydrosystem management,  
22 particularly for understanding and preventing the potential impacts of global change on  
23 ecological functions and services. Macrophyte distribution assessments in rivers are still  
24 primarily realized using field monitoring or manual photo-interpretation of aerial images.  
25 Considering the lack of applications in fluvial environments, developing operational, low-cost  
26 and less time-consuming tools able to automatically map and monitor submerged macrophyte  
27 distribution is therefore crucial to support effective management programs. In this study, the  
28 suitability of very fine-scale resolution (50 cm) multispectral Pléiades satellite imagery to  
29 estimate submerged macrophyte cover, at the scale of a 1 km river section, was investigated.  
30 The performance of nonparametric regression methods (based on two reliable and well-known  
31 machine learning algorithms for remote sensing applications, Random Forest and Support  
32 Vector Regression) were compared for several spectral datasets, testing the relevance of 4  
33 spectral bands (red, green, blue and near-infrared) and two vegetation indices (the Normalized  
34 Difference Vegetation Index, NDVI, and the Green-Red Vegetation Index, GRVI), and for  
35 several field sampling configurations. Both machine learning algorithms applied to a Pléiades  
36 image were able to reasonably well predict macrophyte cover in river ecosystems with  
37 promising performance metrics ( $R^2$  above 0.7 and RMSE around 20 %). The Random Forest  
38 algorithm combined to the 4 spectral bands from Pléiades image was the most efficient,  
39 particularly for extreme cover values (0 % and 100 %). Our study also demonstrated that a  
40 larger number of fine-scale field sampling entities clearly involved better cover predictions than  
41 a smaller number of larger sampling entities.

42

43 **Keywords:** Aquatic vegetation; Remote sensing; Machine learning; Fluvial ecosystem;  
44 Random Forest; Support Vector Regression

## 45 **1. Introduction**

46 The essential role of macrophytes in freshwater ecosystems has been well demonstrated. They  
47 influence the physical, chemical and biological structures of hydrosystems and provide multiple  
48 ecosystem functions and services, such as water quality improvement, stabilization of  
49 streambeds and habitat provision (Carpenter and Lodge, 1986; Dennison *et al.*, 1993; Jeppesen  
50 *et al.*, 1998; Bornette and Puijalon, 2011; Choi *et al.*, 2014). However, their excessive  
51 development can have negative impacts on hydrosystem functioning, for instance, light  
52 attenuation, anoxia, reduction of flow velocity and increase of sedimentation (*e.g.* Caraco and  
53 Cole, 2002; Hussner *et al.*, 2017; Kagami *et al.*, 2019), modifying biotic interactions and  
54 disrupting community assembly (Santos *et al.*, 2011). Submerged species, especially, can also  
55 cause recurring problems for users and managers, *e.g.* inconvenience to water activities,  
56 olfactory nuisances and clogging of water intakes in power plants in case of uprooting (Jadhav  
57 and Buchberger, 1995; Bunn *et al.*, 1998; Sand-Jensen and Pedersen, 1999; Stephan and  
58 Gutknecht, 2002; Martin, 2017) which are still difficult to fully anticipate on large rivers.  
59 Although it is relatively well known that the patterns of spatial and temporal distribution of  
60 aquatic macrophytes result from the interactions between many environmental factors  
61 (hydrology, water temperature, light, nutrients, substrate, grazing), using these factors in  
62 statistical models remains insufficient to predict their distribution within large geographic areas  
63 as fluvial environments. Data on the current distribution of submerged macrophytes in the field  
64 are therefore pivotal, either for direct monitoring, as required by the EU Water Framework  
65 Directive to improve the ecological quality assessment of inland waters (WFD European  
66 Commission, 2000), but also for developing new distribution models.  
67 Mapping and monitoring vegetation distribution are important technical tasks in sustainable  
68 management. Accordingly, numerous monitoring programs have focused on acquiring spatial  
69 information about species composition, maximum depth colonization, density, cover (*i.e.*

70 percentage of the horizontal surface occupied by vegetation), biomass and plant height (Johnson  
71 and Newman, 2011). Those programs have been reviewed in Stocks *et al.* (2019).

72 In fluvial environment, collecting data on submerged aquatic vegetation (SAV) that will  
73 sufficiently represent spatial variation along river reach is difficult, and requires labour-  
74 intensive, time-consuming and sometimes destructive fieldwork. Thus, SAV sampling will be  
75 more and more replaced by indirect mapping methods, especially thanks to remote sensing  
76 tools. Nowadays airborne or satellite sensors provide many observation opportunities at large  
77 scales at a given time, with relatively high spatial and temporal resolutions which are constantly  
78 improving. Actually, multispectral satellite data have been widely used to map the distribution  
79 of macrophytes over large areas (Gullström *et al.*, 2006; Nelson *et al.*, 2006; Dogan *et al.*, 2009;  
80 Tian *et al.*, 2010) and assess the spatio-temporal dynamics of aquatic vegetation (MacAlister  
81 and Mahaxay, 2009; Zhao *et al.*, 2012, 2013). However, many of these studies have focused on  
82 emergent or floating aquatic vegetation due to easier distinction of macrophyte spectral signal  
83 from water. Hyperspectral data have been less considered due to their limitations in terms of  
84 cost, availability, processing and high dimensionality of spectral data (Plaza *et al.*, 2009; Adam  
85 *et al.*, 2010; Mutanga *et al.*, 2012).

86 Recently, significant progress has been made in image processing for repetitive and automatic  
87 submerged macrophyte mapping over large areas combining punctual field data and various  
88 modelling methods. Some researchers have clearly demonstrated the feasibility of submerged  
89 macrophyte mapping using powerful machine learning algorithms (*e.g.* Artificial Neural  
90 Networks, Random Forest, Support Vector Machines or *K*-Nearest Neighbors) (*e.g.* Dogan *et al.*  
91 *et al.*, 2009; Kotta *et al.*, 2013). The main characteristic of these supervised algorithms is to train  
92 a model on a part of the data and test the fitted model on the other part. Compared with other  
93 nonparametric methods, they also have no limitation for the number of independent variables  
94 (*i.e.* adapted to high dimensional data) and do not require normally-distributed

95 variables. Most of these works have been conducted in marine environments, or were limited  
96 to large river basins, wetlands or lakes with the best satellite resolution limited to 2.41 m and  
97 various attempts have been made to resolve issues regarding submerged vegetation (Hedley *et*  
98 *al.*, 2012). However, the insufficient spatial resolution of image data, the spatial variability of  
99 depth and the strong attenuation of light in water are still limitations for remotely mapping SAV  
100 in fluvial environments (Marcus and Fonstad, 2008). To our knowledge, only one study has  
101 focused on submerged macrophytes in rivers combining ultra-light aircraft images to machine  
102 learning algorithms but the obtained map was limited to the presence/absence of canopy  
103 meadows (Durand *et al.*, 2016). Finally, a scientific issue is to develop machine learning models  
104 on high resolution satellite images with high potential in rivers to extend satellite remote  
105 sensing.

106 Additionally, human and financial resources allocated to acquire *in situ* aquatic vegetation data  
107 are generally limited. There is no standard sampling protocol while the quality of macrophyte  
108 cover prediction depends on the sampling strategy adopted. However, to our knowledge, there  
109 is no study in fluvial environments allowing the optimization of SAV monitoring methods, both  
110 in terms of sampling protocol (*i.e.* determining the spatial scale and the minimal number of  
111 sampling plots required) and prediction model choice to generate accurate and continuous map  
112 of riverine submerged macrophytes.

113 In that respect, a new mapping method was here investigated to monitor the distribution of  
114 submerged vegetation in fluvial environments at a very fine-scale resolution (50 cm) while  
115 limiting logistical, financial and human costs. The main objective of this study was to develop  
116 a cover prediction model combining machine learning algorithm, imagery spectral features and  
117 punctual field samples. In that respect, we investigated the performance of two machine  
118 learning regression models (Random Forest and Support Vector Regression) applied to high-  
119 resolution multispectral Pléiades satellite imagery, for automatically mapping macrophyte

120 cover at the scale of a 1 km river section. Both algorithms are reliable and well-known in remote  
121 sensing; they proved to achieve satisfactory results for various and numerous remote sensing  
122 applications in ecology (e.g. Cutler *et al.*, 2007; Hunter *et al.*, 2010; Husson *et al.*, 2017;  
123 Traganos *et al.*, 2018; Villa *et al.*, 2018; Zafari *et al.*, 2019; Sabat-Tomala *et al.*, 2020). Support  
124 Vector Regression is more robust than Random Forest for a lower sample size but the latter is  
125 faster to compute (Mountrakis *et al.*, 2011; Belgiu and Drăguț, 2016; Gholami and Fakhari,  
126 2017).

127 In addition, we discussed several SAV cover sampling strategies involving various numbers of  
128 sampling entities with different sizes in order to guide managers optimizing their monitoring  
129 method. We assessed whether, for a given sampling effort, it is better to use a larger number of  
130 smaller sampling entities or a smaller number of larger entities for SAV cover estimation.

## 131 **2. Materials and methods**

### 132 **2.1. Study area**

133 The study was carried out during September 2017 in the Garonne River, a southwest French  
134 shallow river, approximately 15 km north of Toulouse (43°41'51"N, 1°22'09"E), next to the  
135 city of Seilh (Figure 1a).

136 The study area was about 1 km long and 110 m wide and it is a typical example of the mid-  
137 Garonne ecosystems with abundant macrophyte meadows which develop at low water depth,  
138 from the end of March to early October, and which are mainly constituted of submerged species,  
139 including two dominant species, such as *Myriophyllum spicatum* (Eurasian Water Milfoil, L.  
140 1753) and *Ranunculus fluitans* (River Water-crowfoot, Lam. 1779). A dozen of aquatic  
141 macrophyte species occur at this site and are observed between 0.15 and 2 m depth.

142 The study area is composed of a shallow backwater downstream on the right bank and it is  
143 bordered by a pebble bed on its left bank. A natural weir marks the upstream limit of the site.  
144 Bedrocks, gravels and pebbles are the main substrates and different mixed sediments can be

145 found on half of the study area. Previous *in situ* topography measurements during separate  
146 fieldwork campaigns (combining bathymetric data collected with a single beam echosounder  
147 and elevation measurements of the riverbed, banks and overflow areas) and 2D hydrodynamics  
148 simulations were used to determine water levels at the whole-site scale during the macrophyte  
149 sampling period (§2.2.). Water levels were relatively low, with average depths varying between  
150 0.05 and 1.30 m over more than half of the site during this period. The downstream zone was  
151 deeper varying mainly between 1.30 and 3.50 m, and with a maximum depth of 4.05 m on the  
152 right side of the central channel.

## 153 **2.2. Field data collection**

154 Field survey of submerged macrophytes was conducted during the second half of September  
155 2017 to get observed data of macrophyte covers at a period with available Pléiades images and  
156 a non-turbid water column. Numerical simulations based on a SAV model indicated that the  
157 biomass of macrophytes on this site and at that date was still very close to the peak biomass  
158 that occurred during the second half of August (unpublished results). A total of 55 sampling  
159 plots of 9 m<sup>2</sup> (*i.e.* 3 m side PVC square frames) were distributed all over the study site (Figure  
160 2a). The sampling protocol focused on homogeneous areas of macrophyte meadows and was  
161 oriented in order to sample several combinations of substrate, depth and meadow abundance  
162 classes; it also included numerous open water plots (Table 1). Each plot (P) was divided into  
163 16 quadrats (Q) of 0.75 m side (Figure 2b). Then the total cover (*i.e.* referring to the cover of  
164 all species) of SAV within each quadrat was estimated by two subaquatic observers. A total of  
165 880 quadrats were thus sampled individually. Finally, we averaged the total cover from the 16  
166 quadrats of each plot to get the total cover at the plot scale. The purpose here was to compare  
167 two field sampling levels: the plot level (*i.e.* 55 plots) versus the quadrat level (*i.e.* 880  
168 quadrats).

169 The centre of each sampling plot was geolocated using a portable real time kinematic global  
170 positioning system (RTK-GPS) receiver (EMLID Reach RS™) with +/- 3 cm accuracy. The  
171 RTK-GPS reference receiver was located on field, close to the right river bank of the study area  
172 (*i.e.* <1 km from all sampling plots) and free of obstacles to ensure maximum exposure to  
173 radiometric signals. Both RTK-GPS receivers used GNSS signals from GPS and GLONASS  
174 satellites. All plots were North oriented. The geographical information system QGIS 2.18 was  
175 used to digitize the field data.

### 176 **2.3. Remote sensing data acquisition into the Garonne River**

177 Two types of optical measuring instruments were used to acquire remote sensing data:  
178 unmanned aerial vehicles (UAV) and satellite with high spatial resolutions. UAV imagery  
179 provided an overview of the distribution of submerged macrophytes meadows at the site scale  
180 whereas satellite image was used for the development of the automatic mapping method.

#### 181 **2.3.1. UAV imagery**

182 Aerial photos were taken mid-September 2017 with a Phantom 4 Pro, a drone developed by  
183 DJI™ (DJI™, Shenzhen, China). We used a flying height of 30 m. The along- and across-track  
184 image covers were set to 70-80 %. The study site was surveyed in three flight blocks of 10-15  
185 min (Figure 3a). Flight missions were programmed with the Litchi application software  
186 available for Android devices. Flight weather conditions were sunny with light wind at ground  
187 level. The camera used in this work was a three-band (red, green and blue) DJI™ digital camera,  
188 equipped with a 1" CMOS / 20-megapixel sensor camera, with a focal length of 24 mm, and a  
189 F-number of 2.8, providing an image size of 5472 × 3648 (columns × rows) (Figure 3b).

190 Mosaicking was processed using the commercial software Agisoft© Metashape Professional  
191 Edition (Saint Petersburg, Russia) with a spatial resolution of 1 cm (Figure 3c). Georeferencing  
192 was performed on QGIS 2.18 using seven ground control points which were taken *in situ* with  
193 the RTK-GPS with a linear transformation and nearest neighbors resampling. Then a map of

194 the total cover of submerged macrophytes was created on QGIS 2.18 by visual interpretation  
195 of the orthomosaic at each 3×3 m cell of an overlaid grid. This photo-interpreted cover map  
196 provided a field reference of the study site.

### 197 **2.3.2. Satellite imagery**

198 One Pléiades pan-sharpened image of 100 km<sup>2</sup> surrounding our study site (Figure 1b) was  
199 acquired on September 27, 2017, thanks to the “Initiative for Space Innovative Standards”  
200 (ISIS) program that results from a cooperation between Airbus Defence and Space and the  
201 French “Centre National d’Etudes Spatiales” (CNES). Pan-sharpening, resulting from a fusion  
202 process, corresponds to a multispectral image including 4 spectral bands from the visible (Red:  
203 0.59-0.71 μm, Green: 0.50-0.62 μm, Blue: 0.43-0.55 μm or RGB) through the near-infrared  
204 (NIR: 0.74-0.94 μm) with the spatial resolution of panchromatic images (*i.e.* 50 cm) obtained  
205 by the Pléiades-1A satellites. The ortho-image was cloud free with minimal glint and corrected  
206 from acquisition and terrain off-nadir effects by the providers.

207 The raw Pléiades image was already ortho-rectified before delivery (with 2.5 m accuracy  
208 according to Grazzini *et al.* (2013). Its georeferencing has been improved using a linear  
209 transformation and nearest neighbors resampling. This process required 17 ground control  
210 points for which 3 cm accuracy GPS positions were determined with the RTK-GPS. This  
211 process allowed a new planimetric accuracy of the Pléiades image of *ca.* 0.35 m.

212 Then, radiometric corrections have included conversion of pixel digital number values  
213 (encoded on 16 bits) to Top of Atmosphere (TOA) reflectance for each spectral band (*i.e.* RGB  
214 and NIR) in order to normalize each spectral band in a a continuous range between 0 and 1.  
215 This process was made using the “Geosud TOA Reflectance” plugin available on QGIS 2.18.  
216 Finally, a vector mask of the riparian area and dewatered banks was used to produce a minor  
217 riverbed-only image.

218

## 219 **2.4. Modelling macrophyte cover using satellite imagery**

220 Modelling the relationship between SAV cover and its spectral reflectance is essential for  
221 automatically mapping macrophytes. We used machine learning regression algorithms to  
222 predict macrophyte cover from high-resolution reflectance data based on Pléiades image.  
223 Figure 4 summarizes the main steps of our approach.

224 The predictive model was developed using free geomatics and statistics tools, such as QGIS  
225 2.18 spatial analysis technologies to pre-process the remote sensing data and several R libraries,  
226 such as “rgdal”, “randomForest” and “e1071”, to develop the regression models and to generate  
227 the predicted distribution maps at the spatial resolution of the Pléiades images (*i.e.* 0.5 m of  
228 resolution). Finally, their resolution had been degraded to 3 m resolution for comparison with  
229 the photo-interpreted cover map (*i.e.* 3 m).

### 230 **2.4.1. Defining predictive variables among spectral data**

231 The use of red and especially of NIR band for submerged macrophyte detection is still debated  
232 because of their relatively fast absorption in water (Fyfe, 2003; Heege *et al.*, 2004; Dogan *et*  
233 *al.*, 2009; Chen *et al.*, 2018b). However, many submerged macrophytes reach the surface during  
234 low water period by creating a canopy having stronger contribution to the signal, potentially  
235 making these two bands suitable for the detection of aquatic vegetation. In addition, the use of  
236 normalized spectral indices could improve prediction models (Bradley and Fleishman, 2008).  
237 Different datasets, derived from the pre-processed satellite image, were therefore used to assess  
238 the potential of multispectral bands and vegetation indices to detect submerged macrophytes,  
239 and to examine the effect of spectral features on the quality of cover prediction. The four  
240 spectral reflectance datasets were the following:

- 241 (i) The first dataset was obtained from the Pléiades concatenated spectral data from the  
242 visible bands (red, green and blue bands (RGB)). This dataset was referred to as  
243 “dataset 0” (3 variables);

- 244 (ii) The second dataset (“dataset 1”) combined RGB and near-infrared (NIR) bands (4  
245 variables);
- 246 (iii) The third dataset (“dataset 2”) included two widely-used vegetation indices  
247 reflecting the characteristic spectral signature of active vegetation: the Normalized  
248 Difference Vegetation Index,  $NDVI=(NIR-R)/(NIR+R)$ , and the Green-Red  
249 Vegetation Index,  $GRVI=(G-R)/(G+R)$  (2 variables). According to Cho *et al.*,  
250 (2008) NDVI could only be useful in shallow clear waters (<0.5 m depth) because  
251 of fast absorption of the NIR band while GRVI, using red and green bands, decrease  
252 less rapidly depending on the depth and could perform better in detecting changes  
253 in canopy vegetation (Motohka *et al.*, 2010; Chen *et al.*, 2018a);
- 254 (iv) The last dataset (“dataset 3”) included the RGB and NIR bands and the two GRVI  
255 and NDVI indices (6 variables).

#### 256 **2.4.2. Extraction of reflectance data based on sampling entities**

257 Our field database was composed of 55 plots or 880 quadrats (each of them being characterized  
258 by a value of total macrophyte cover expressed in percentage). Each sampling entity (plot or  
259 quadrat) was composed of a given number  $n$  of pixels from the Pléiades image, each pixel being  
260 associated with a reflectance value in several spectral bands (or vegetation indices). The  $n$   
261 values of reflectance of each spectral band are extracted using the *extract()* function from the  
262 R *raster* package. Then an area-weighted average reflectance of the  $n$  pixels belonging to each  
263 sampling entity was computed.

#### 264 **2.4.3. Machine learning regression models and SAV mapping**

265 We tested the performance of two reliable machine learning algorithms to build a regression  
266 model that predicts SAV cover at the scale of the site: the Random Forest (RF) and Support  
267 Vector Regression (SVR) algorithms. For each model, the remote sensing variables (including

268 vegetation indices for “dataset 2” and “dataset 3”) were treated as independent variables and  
269 the total cover was treated as a dependent or response variable. Detailed reviews of SVR and  
270 RF in remote sensing can be found in Mountrakis *et al.* (2011) and Belgiu and Drăguț (2016),  
271 respectively.

272 Regression analysis was carried out on the four datasets and at the two sampling levels (plots  
273 and quadrats). This procedure allows us to assess the effect of the spectral dataset and the  
274 relevance of the two field sampling levels.

275 We then explored the effects of the number and size of the sampling entities on prediction  
276 quality. Firstly, the effect of the sampling scale on cover prediction was addressed by comparing  
277 the results obtained using datasets with 55 entities but of different size: 0.75 m, 1.5 m  
278 (corresponding to a group of 4 quadrats per plot) and 3 m side. Secondly, the number of quadrats  
279 was gradually decreased from the initial dataset (880 quadrats) by randomly selecting 1, 2, 4 or  
280 8 quadrat(s) per plot. This random selection of quadrats generated configurations with 55, 110,  
281 220 or 440 quadrats. In addition, a subset of the plots was used to get a configuration with 15  
282 and 30 quadrats. Thanks to an oriented random selection function, we ensured that the quadrat  
283 selection maintain a similar distribution of macrophyte abundance classes, as that of the initial  
284 880 quadrats dataset (Table 1). The purpose here was to define the minimal number of sampling  
285 entities that allowed a satisfactory result. This was achieved by a learning curve analysis.

286 In total 72 models (combining 2 algorithms, 4 spectral datasets and 9 field sampling  
287 configurations) of cover prediction were run and compared with each other in order to  
288 determine the best one. The key steps of the method are detailed below.

289 • ***Tuning, training and testing the models***

290 In the case of a limited number of samples, the *k*-fold cross-validation technique is recognized  
291 as a valuable approach to split samples into *k* subsets of roughly equal size (Hastie *et al.*, 2009).

292 For a given combination of hyperparameters,  $k-1$  partitions are used iteratively ( $k$  times) to train  
293 the model and to test it on the remaining partition. Thus, it potentially allows each sample to be  
294 used  $k$  times for multiple training or testing, with the purposes of (i) improving the learning  
295 process (fine tuning), (ii) using independent datasets for training and testing, and (iii) limiting  
296 overfitting (Anguita *et al.*, 2012; Ramezan *et al.*, 2019). The final model performance is  
297 calculated by averaging the  $k$  computed errors for the selected hyperparameters. Finally, the  
298 best prediction model corresponds to the model built with hyperparameters which generate the  
299 highest validation score (*i.e.* the lowest generalization error or test error). Additionally, a  
300 commonly-used method to reduce the variability in chosen parameters and the standard  
301 deviation of performance estimates of the tuned model, is to run a repeated  $k$ -fold cross-  
302 validation, named  $J$ - $k$ -fold cross-validation (Moss *et al.*, 2018). We therefore used in this study  
303 a repeated  $k$ -fold cross-validation, *i.e.* with  $k$  fixed to 10, as is commonly chosen in the literature  
304 (Jung, 2017) and we set  $J$  to 10. Thus, in each round, a number of sampling entities was  
305 randomly selected, with 90 % used as training data and 10 % as test data.

306 During the learning step, the different combinations of hyperparameter values are examined  
307 and tuned for the calibration of the model. The combination of values that generate the lowest  
308 error on the test set is assumed to be optimal (Brenning, 2012; Cracknell and Reading, 2014;  
309 Sharma *et al.*, 2017). For the RF algorithm, the number of regression trees  $n_{tree}$  ranged from  
310 25 to 1000 (with a step of 25) to test the sensitivity of this parameter. The  $n_{tree}$  values that  
311 yielded the lowest error were selected for each dataset. Due to conflicting reports in the  
312 literature concerning the potential influence of the  $mtry$  parameter (*i.e.* the number of predictive  
313 variables) on prediction performance (Cutler *et al.*, 2007; Strobl *et al.*, 2008), the  $mtry$  value  
314 was tested from 1 to the maximum number of spectral bands for each dataset (*e.g.* maximum of  
315 6 for “dataset 3”). Concerning the SVR algorithm, the commonly used gaussian radial basis  
316 kernel function (RBF) was applied, given its traditional superior performance compared to other

317 kernels (*i.e.* linear or polynomial) (Kavzoglu and Colkesen, 2009). The different parameters  
318 ranges were as follows: the regularization parameter  $C$  (*i.e.* cost) ranged from 0.1 to 1000 (by  
319 a factor of 10),  $\epsilon$  was fixed to 0.1 and the width of the RBF kernel function  $\gamma$  ranged from  $2^{-5}$   
320 to  $2^5$  (by a factor of 2). Predictive variables were standardized.

### 321 • **Model evaluation**

322 Two statistics were employed to evaluate the quality of the model predictions:

- 323 - The coefficient of determination ( $R^2$ ) which varies between 0 and 1, to account for the  
324 goodness-of-fit between observations and predictions, and to define how much variance is  
325 explained by the model.
- 326 - The root-mean squared error (RMSE) to assess the predictive power of the model.

327 These statistics were computed at the scale of the field entities, as well as the scale of the site,  
328 using punctual visually-estimated covers for the former (depending on the configuration  
329 tested, for instance 55 plots or 880 quadrats), and the cover map obtained from drone imagery  
330 photointerpretation for the latter. However, only the metrics computed at the site scale will be  
331 discussed here to address the predictive quality of our models.

332 Difference maps between predicted and observed covers were also computed to pinpoint local  
333 prediction errors (expressed as cover percentages; a difference below or above 0 indicating an  
334 underprediction or an overprediction, respectively).

## 335 **3. Results**

### 336 **3.1. Observed macrophyte distribution on the Garonne River**

337 Monospecific and plurispecific meadows were observed in the field. Sampling entities with  
338 high cover values generally included the two dominant species (see §2.1) while the remaining  
339 species were observed at lower densities, except for *Potamogeton nodosus* and the *Elodea*  
340 species which were locally abundant, notably in the backwater. Species were distributed

341 according to their ecological requirements (flow, turbidity). Table 1 summarizes the distribution  
342 of densities according to the adopted field sampling strategy. Table S1 (see supporting  
343 information) provides the database from field sampling.

344 The photointerpretation of the drone mosaic (*i.e.* the orthomosaic) (Figure 5a) highlighted the  
345 spatial distribution of macrophytes and revealed variability in the distribution of macrophyte  
346 meadows. These were found mainly upstream and halfway along riverbanks, but also at the  
347 backwater on the right bank. Overall, higher covers were found along the left bank, in areas of  
348 shallow depths (0.1-1 m) especially on mixed substrates. Lower densities were observed in  
349 deeper areas and/or on rock bed. Downstream meadows seemed to be patchier along the left  
350 bank. Rare meadows were present in the centre of the channel. Bare areas (*i.e.* with 0 % cover)  
351 were the most represented on the site, followed by areas with covers between 1 % and 10 %,  
352 especially present along the right bank, while areas with covers between 11-25 % and above 75  
353 % were mostly located along the left bank. Based on this cover distribution and on point  
354 biomass measurements at given sampling plots (data not shown), the total biomass on the site  
355 was estimated to be 2.6 t of dry matter for a surface of *ca.* 10 ha.

## 356 **3.2. Predicted cover of submerged macrophytes**

357 All the machine learning results were analysed in terms of statistical performance and  
358 prediction quality (difference between predicted and observed covers). For each regression  
359 model, the  $R^2$  and RMSE statistical results, as well as the range of the predicted cover values,  
360 are provided in the supporting information– Table S2. Figures S1 to S4 also group together the  
361 whole map results in the supporting information. Here we focus on the most relevant results.

### 362 **3.2.1. Effect of the spectral features**

363 Analysing the effect of spectral bands on macrophyte cover prediction has shown meaningful  
364 differences between the statistical metrics ( $R^2$  and RMSE) of the different datasets tested,  
365 regardless of the sampling level (*i.e.* plot or quadrat) considered.

366 Prediction models built on “dataset 2” (*i.e.* NDVI and GRVI) obtained the least satisfactory  
367 mapping and statistical results, regardless of the algorithm or the sampling level. The  
368 determination coefficients ( $R^2$ ) (*i.e.* similarity levels between predicted and photo-interpreted  
369 covers) were of 0.4-0.5 and the prediction errors (RMSE) ranged between 25 % and 27 % (Table  
370 S2). Overall, the bare areas (*i.e.* open water) were poorly predicted with important over-  
371 prediction, particularly with the Support Vector Regression (SVR) algorithm. Cover maps were  
372 improved to some extent with the Random Forest (RF) algorithm: boundaries of macrophyte  
373 patches were more clearly distinguished (Figures S1-S2). Moreover, the difference cover maps  
374 showed under-predicted areas, particularly for the meadows located along the left and right  
375 banks (Figures S3-S4).

376 In contrast, no meaningful difference with respect to the evaluation metrics, were found for the  
377 models based on the three remaining datasets and using either RF or SVR. All of them usually  
378 indicated high goodness-of-fit at the study scale ( $R^2$ ), between 0.62 and 0.75, as well as a good  
379 prediction power with a RMSE of 21 % on average (maximum of 24 %) (Table S2).

380 However, regarding RF, a poorer prediction of the bare areas was noticed in case of “dataset  
381 0”, while the results obtained with datasets 1 and 3 were equivalent (see Figures S1, S3 and  
382 Table S2 for details). Adding NIR improved prediction quality in the bare areas, particularly at  
383 the quadrat sampling level.

384 For SVR, only the maps obtained with datasets 0 and 1 represented relatively well the spatial  
385 distribution of macrophyte meadows on the Garonne River (Figure S2). Indeed, with “dataset  
386 3”, high densities were better predicted (few under-prediction errors) than with “dataset 2” but  
387 bare areas still presented considerable over-prediction errors (differences up to 28 % between  
388 predicted and observed covers) (Figure S4). Furthermore, the predictive advantage of adding  
389 NIR (“dataset 1”) was variable depending on the 2 sampling levels (*i.e.* plots or quadrats). For  
390 instance, improved results were noticed for models based on “dataset 0” (involving the 3 bands

391 in the visible) and on quadrats: this dataset yielded the lowest difference between predicted and  
392 observed covers (Figure S4).

393 For the rest of result analysis, we excluded datasets 2 and 3 since they produce models that were  
394 poorly predictive, and we focused on models based either on RF and involving “dataset 1” or  
395 on SVR and involving datasets 0 and 1.

### 396 **3.2.2. Effect of the learning algorithm**

397 The two machine learning algorithms tested here and used on our two field datasets (55 plots  
398 and 880 quadrats) were able to predict the distribution and macrophyte cover on the Garonne  
399 River, reaching a maximum predicted cover value of 100 % (Table 2). Whether for plots or  
400 quadrats, no statistical pattern was highlighted through the comparison of the two algorithms.  
401 Regression models showed high  $R^2$ , oscillating between 0.67 and 0.74 and low prediction errors  
402 (RMSE around 19-21 %), even if SVR models reached slightly higher performance (Table 2).

403 Overall the predicted cover maps reproduced the observed spatial distribution of submerged  
404 macrophytes, with dense meadows along the left bank over more than half of the site (Figure  
405 5c and Figures S1-S2). Nevertheless, there were two areas where the cover prediction error  
406 remained high (differences up to 100 % between the predicted and observed covers), namely in  
407 the backwater area, and downstream, within the dense meadow located along the left bank,  
408 where high densities (> 75 %) were detected as bare areas. These differences were identified  
409 by the dark red areas in the difference maps (*e.g.* Figure 5d).

410 Besides, detecting the patchier distributions downstream, within the backwater and along the  
411 right bank seems to be more or less difficult depending on the algorithm. In that respect, the  
412 prediction accuracy of macrophyte cover along the riverbank in the middle of the study site was  
413 lower using the SVR algorithm than using the RF algorithm (Figures S1-S2).

414 In addition, a visual comparison between observed and predicted cover data allowed to point at  
415 certain characteristics of the regression algorithms (see Figures S3-S5 in supporting information  
416 for details). If we focus on results based on SVR, at the entity scale, data were globally noisier  
417 especially with important discrepancies for extreme values of cover (*i.e.* 0 and 100 %) (Figure  
418 S5). In that respect, more prediction errors were observed at these cover densities at the site  
419 scale, in particular for meadows upstream and within the central channel (Figure S4). In  
420 addition, average densities (*i.e.* those between 25 and 75 % of cover) were hardly predicted  
421 with SVR models. Indeed, those areas were often underestimated by this algorithm, with  
422 differences between predicted and observed cover reaching a maximum of 50 %. On the other  
423 hand, the fit of the regression model with the observed covers was better using RF, compared  
424 to SVR (Figure 5b and Figure S5). The different covers seemed globally well predicted by RF,  
425 with the different classes of macrophyte abundance better discriminated, even if the highest  
426 covers (*i.e.* above 90 %) were often slightly under-predicted (maximum of difference under 28  
427 %). Finally, the bare areas were better represented by this algorithm and, in case of over-  
428 prediction, the maximum difference was still limited to 20 %, particularly along the riverbeds  
429 (Figure 5d).

430 Consequently, the RF algorithm when associated with “dataset 1” was considered as the most  
431 convincing one for predicting the spatial distribution of submerged macrophyte cover.

### 432 **3.2.3. Effect of field sampling strategy: 55 plots vs 880 quadrats**

433 The comparison of the two sampling levels (plots vs. quadrats) highlighted clear differences  
434 between algorithms in their abilities to detect macrophyte patches. Whatever the algorithm,  
435 goodness-of-fits and spatial errors were not notably different between the two sampling entities  
436 although a model using 880 quadrats seemed to show performances somewhat higher than those  
437 obtained with 55 plots dataset (Table 2). Even if both of these sampling levels predicted a wide  
438 range of cover percentages, regression models based on quadrats predicted better the largest

439 cover values, especially those above 98 % (Table 2). Furthermore, the limits of the meadow  
440 patches seemed less well-defined with a model using 55 plots, compared to 880 quadrats.  
441 Finally, the difference maps revealed that the differences between predicted and observed  
442 covers were reduced using 880 quadrats (differences less than 20 % on average) (Figure S3).  
443 Consequently, although models based on the two field sampling levels were able to distinguish  
444 the spatial distribution of meadows, considering 880 quadrats instead of 55 plots showed better  
445 results. Overall, the best model for macrophyte cover prediction was obtained with RF using  
446 NIR band (§3.4.1 and 3.4.2) and based on field data sampled within 880 quadrats (Figure 5c).  
447 However, it remains to be seen whether these better performances are due to a finer size and/or  
448 to more numerous entities.

#### 449 **3.2.4. Effect of the size of sampling entity**

450 The comparison of the models based on 3 different sizes (3 m, 1.5 m and 0.75 m side) of the  
451 sampling entities and with a fixed number of entities ( $n = 55$ ) showed very clear differences  
452 regarding cover prediction accuracy.

453 Actually, even if no meaningful difference in model metrics ( $R^2$  and RMSE) could be seen  
454 between the different scales, the RF model built on sampling entities of 3 m side presented  
455 slightly better  $R^2$  and RMSE compared to models using smaller sampling entities ( $R^2 = 0.70$ ,  
456 RMSE = 21.3 %) (Table S2). Boundaries between the different macrophyte patches were better  
457 defined too (Figure S1). The difference maps also showed the least spatial difference between  
458 predicted and observed covers, with average differences amounting to +8 % in bare areas and -  
459 40 % within macrophytes meadows. Spatial differences locally reached +12 % in bare areas  
460 and -72 % within meadows (Figure S3). Models based on quadrats of 0.75 m side showed the  
461 worst performances ( $R^2 = 0.62$ , RMSE = 23.7 %) (Table S2) and poor cover map results (with  
462 local prediction differences reaching +20 % in bare areas and -90 % within meadows) (Figure  
463 S3). With sampling entities of 1.5 m side, results were improved but still worse than a model

464 based on 3 m side sampling entities. Similar results were also observed with the SVR algorithm  
465 (see Figures S2, S4 for details).

466 Therefore, for a limited number of entities ( $n = 55$ ), predicting cover using 3 m side entities  
467 seemed more appropriate when addressing the spatial patterns of macrophytes. However as  
468 shown above, this sampling strategy was not fully effective for accurate prediction of cover, in  
469 particular compared to a sampling strategy with more entities.

### 470 **3.2.5. Effect of the number of entities**

471 Analysis of learning curves, which described the evaluation metrics of models based on  
472 different numbers of quadrats, showed that the maximum performance was reached for 110  
473 quadrats (and above) whatever the spectral dataset and algorithm used (Figure 6 and see Figure  
474 S6 in supporting information for the SVR results). More precisely, the  $R^2$  and RMSE remained  
475 stable from 110 quadrats for the RF models (Figure 6). Regarding the SVR, the same  
476 observations could be made, though the metrics were really stable from 220 quadrats for  
477 “dataset 1” (RGB NIR).

478 In terms of sampled surface, 220 quadrats correspond to 55 entities of 1.5 m side (*i.e.* 4 grouped  
479 quadrats of 0.75 m side each) and a 220 quadrats-based model has shown to be more performing  
480 than a 55 plots-based one (and, a fortiori, than a model based on 55 entities of 1.5 m) (Figure 6  
481 and Figure S6). Consequently, for a same sampled surface, cover estimate will be better with  
482 models involving more but smaller entities than with models using a limited number of larger  
483 entities.

## 484 **4. Discussion**

485 Surveying submerged macrophytes using remote sensing is somewhat more difficult than  
486 surveying terrestrial vegetation because of water reflectance issues (linked to the strong  
487 reflection of the water surface and attenuation of light by the water column) and the limited

488 spatial resolution of most sensors (Nelson *et al.*, 2006; Underwood *et al.*, 2006). In this study,  
489 we developed a new method for automatically mapping the cover of SAV on a river section  
490 during the peak of biomass. This method included very high-resolution (50 cm) spectral data  
491 from Pléiades pan-sharpened image combined with a machine-learning algorithm and an  
492 optimized method of macrophyte cover sampling.

#### 493 **4.1. Developing a remote sensing method for automatic mapping of vegetation cover**

##### 494 **4.1.1. Choosing an appropriate machine learning algorithm**

495 The achievement of this remote sensing method did not really depend on the two regression  
496 algorithms tested (RF and SVR). Between both of them statistical results were similar, with  
497 only small differences in the predicted cover maps. They showed reasonable fitting capacity  
498 ( $R^2$  around 0.7) and relatively low prediction errors (RMSE < 25 %). In fact, a correct  
499 parameterization of different machine learning algorithms must lead to similar results.  
500 However, fewer local prediction errors were observed for the RF models. The spatial  
501 boundaries of the meadow located along the right bank were also better defined. Actually, SVR  
502 is more sensitive to parameter assignment than RF (among them the choice of the  $\gamma$  parameter  
503 which allows optimal data discrimination) during the tuning step (§2.4.3) that can significantly  
504 alter the performance of SVR algorithms (Brown *et al.*, 1999; Mountrakis *et al.*, 2011).  
505 Additionally, the RF algorithm is known to be effective in quickly handling high dimensional  
506 data and multicollinearity (Belgiu and Drăguț, 2016). It was thus considered as the most  
507 appropriate machine learning algorithm for river macrophyte mapping.

##### 508 **4.1.2. Determining the best spectral features**

509 For models based on the RF algorithm, the best predicted results were obtained using datasets  
510 combining visible spectral and NIR bands (vegetation indices such as NDVI and GRVI were  
511 of no use for cover prediction). Adding NIR to visible spectral bands also seemed to slightly

512 benefit the prediction of low covers but did not systematically improve the quality of the  
513 prediction in the areas of higher expected covers. The advantage of adding NIR depended on  
514 the algorithm. Indeed, the use of NIR and the indices derived from it (NDVI) are still widely  
515 discussed in the aquatic remote sensing literature, because of the large absorption by water of  
516 red and, even more, NIR wavelengths (*e.g.* Pegau *et al.*, 1997; Fyfe, 2003; Cho *et al.*, 2008;  
517 Silva *et al.*, 2008). Moreover, even if the green wavelengths normally provide greater light  
518 penetration in turbid waters, green and red regions are considered as the best ones for  
519 submerged macrophyte sensing (Fyfe, 2003; Han and Rundquist, 2003; Williams *et al.*, 2003;  
520 Pinnel *et al.*, 2005). The use of GRVI did not yet compensate for the low performance of NDVI  
521 in our study. Besides, a study carried out on wetlands has shown that red and NIR regions could  
522 be saturated beyond a certain biomass density (Peñuelas *et al.*, 1993; Mutanga and Skidmore,  
523 2004). Finally, Chen *et al.* (2018b) have clearly demonstrated that NIR should only be used in  
524 the case of very shallow aquatic environments (< 1 m depth). NIR could thus be useful for  
525 predicting macrophyte meadows along the riverbanks, or when submerged macrophytes form  
526 a canopy just below the water surface, which is often the case in eutrophic waters at the biomass  
527 peak period.

#### 528 **4.1.3. Optimizing field sampling**

529 Training sample size also influenced the prediction accuracy of macrophyte cover. With a  
530 limited number of entities ( $n = 55$ ), using 3 m side entities was more effective than using smaller  
531 entities. Indeed, as the geolocation of the Pléiades image was not perfect, it is likely that the  
532 small sampling entities did not exactly match with the satellite image. In this case, increasing  
533 the size of the sampling entity would solve this problem and could explain the better results for  
534 3 m side size entities. However, for the RF (and SVR) regression models based on 55 plots,  
535 local prediction errors were larger, particularly for bare areas, than the model based on 880  
536 quadrats. For a same surface sampled, the latter actually showed the best predictions, even if

537 certain areas still presented local under-prediction errors. In fact, the bigger the training dataset  
538 the better the algorithm will learn. Predicted macrophyte meadows were particularly well  
539 distributed and, when present, local prediction errors were the lowest in comparison with other  
540 field sampling strategies. Both meadows along the riverbanks were well defined. Using a power  
541 relationship ( $R^2 = 0.8$ ) between shoot biomass and cover previously established (unpublished  
542 study, see Figure S7 in supporting information), shoot biomass at the site scale was estimated  
543 to 2.3 t of dry matter for 10 ha. This is relatively consistent with the shoot biomass derived from  
544 photointerpretation (*i.e.* 2.6 t of dry matter) and confirmed the performance of our model.

545 Machine learning algorithms were also influenced by the number of training data and their  
546 distribution into the study site. Learning curves confirmed that up to 55 quadrats (*i.e.* entities  
547 of 0.75 m side) the number of data was insufficient to make valuable predictions. Considering  
548 a higher number of training data, particularly in areas where reflectance values were highly  
549 fluctuating, would improve prediction accuracy by accounting for larger reflectance variability.  
550 Our results showed that for models using 110 (and above) quadrats, predicted cover maps and  
551 model statistics were very similar to those based on 880 quadrats; the estimates with 220  
552 quadrats were clearly improved in comparison to those obtained using 55 entities with higher  
553 sampling surface (1.5 or 3 m side), despite an equal or smaller total sampled surface. This study  
554 revealed the higher performance of sampling numerous (even smaller) entities. This criterion  
555 is particularly interesting regarding submerged macrophyte monitoring as the eye estimation of  
556 vegetation covers is relatively difficult at a 3 m scale, particularly in deep areas, due to the  
557 difficulty to get an overview of macrophyte cover on such a surface.

558 Compared to 55 entities of 3 m side (*i.e.* plots), the sampling surface is also reduced by 8 using  
559 110 entities of 0.75 m side (*i.e.* quadrats). Even if using numerous smaller entities implies  
560 additional time in the field for movements in water and for GPS coordinate acquisition, this  
561 should be largely compensated by the time saved to estimate total cover within entities of

562 smaller size , and by handling a more ergonomic sampling accessory (*i.e.* a smaller PVC frame).  
563 From our experience, the sampling time spent on an entity was almost linearly linked to its  
564 surface, because of the need to split large entities in smaller sub-entities to get accurate cover  
565 estimates. Thus, in comparison with 55 plots, the overall sampling effort should be significantly  
566 reduced when using 110 quadrats, while increasing prediction performance. Consequently, for  
567 future monitoring campaigns along a river section of 1 km long, we recommend estimating  
568 cover in 100 to 200 individual quadrats of a surface comprised between 0.5 and 1 m<sup>2</sup>, which  
569 have shown to be optimal to obtain representative macrophyte mapping using Pléiades imagery.  
570 Despite our method performing reasonably well, macrophyte cover was still underestimated to  
571 some extent (maximum local difference between predicted and observed cover still reaching 28  
572 %) and some bare areas were overestimated (maximum local difference reaching 22 %). Some  
573 studies reviewed in Guo *et al.* (2017) have discussed the effect of imbalanced data on the quality  
574 of machine learning classification. When some field attributes are infrequently present they can  
575 be most likely predicted as rare occurrences, undiscovered or ignored, or assumed as noise or  
576 outliers which results to more prediction errors of certain covers (Ali *et al.*, 2005). However, as  
577 pointed out by Visa and Ralescu (2005) perfect balanced training data is not a guarantee to  
578 improve a classifier performance. Field sampling has to be representative of the study site  
579 (Petersen *et al.*, 2005). Therefore, regardless of the distribution of our macrophyte abundance  
580 classes (Table 1), it is possible that better balancing the entity numbers, particularly for the  
581 extreme covers (*i.e.* 0% and 100 %), will improve mapping results in future investigations.  
582 Indeed, this sampling strategy would include more reflectance variability, as it is particularly  
583 observed in open water entities.

#### 584 **4.2. External factors influencing model quality**

585 Isolating plant signal from the water column interference is still the main challenge of remote  
586 sensing of SAV, due to the low contrast (Williams *et al.*, 2003) and to the inherent difficulties

587 in interpreting reflectance values of water (Peñuelas *et al.*, 1993; Lehmann and Lachavanne,  
588 1997). Numerous studies have revealed that the spectral signal of SAV can also be limited by  
589 environmental and biological factors, such as water depth, turbidity/transparency, distance  
590 between vegetation canopies and water surface (Maritorena *et al.*, 1994; Han and Rundquist,  
591 2003; Vis *et al.*, 2003; Dogan *et al.*, 2009; Liew and Chang, 2012). For instance, local  
592 overestimations of low covers were observed in the central channel or downstream of our study  
593 site, where water flow and depth are high or where suspended matter is highly concentrated  
594 with a thick layer of mud on the bottom. For future investigations it would be interesting to  
595 determine if predictions of macrophyte cover could be improved by including substrate types  
596 or measures of water clarity (*e.g.* Secchi depth, chlorophyll a content) in our models, as stated  
597 by Nelson *et al.* (2006).

598 Some of the 12 submerged species in our study site with low height (*e.g.* *Elodea canadensis*)  
599 could also be undetected. Several studies about submerged vegetation mapping have shown that  
600 non-canopy forming aquatic vegetation species generally lead to more detection errors (Vis *et*  
601 *al.*, 2003; Valta-Hulkkonen *et al.*, 2005; Wolter *et al.*, 2005). It has been reported that SAV can  
602 be remotely sensed to a maximum depth between 2 m and 3 m (Han and Rundquist, 2003;  
603 Sawaya *et al.*, 2003). Finally, low concentrations of some photosynthetic pigments in plant  
604 leaves, such as chlorophylls a and b, carotene and xanthophylls could also affect the spectral  
605 reflectance among vegetation (Kumar *et al.*, 2001).

## 606 **5. Conclusion**

607 Our results provided further evidence that macrophyte cover can be reasonably well predicted  
608 with automated regression procedures based on machine learning algorithms and a limited  
609 number of sampling entities.

610 Remote sensing of riverine submerged macrophytes by pansharpned Pléiades imagery  
611 associated to a Random Forest algorithm appeared to be a viable and valuable tool for

612 estimating biophysical measures, such as macrophyte cover, at very high spatial resolution (50  
613 cm) on a 1 km site on the Garonne River. Performance metrics were promising with  $R^2$  above  
614 0.7 and prediction error rates around 20 %. In this paper we provided a new, efficient and less  
615 time-consuming tool for monitoring SAV which should help steering environmental  
616 management actions such as SAV restoration projects or overgrowth management.  
617 There is a significant opportunity for applying such a promising method to the multi-date  
618 monitoring of SAV in freshwater river environments. Indeed, the monitoring and mapping of  
619 macrophyte meadows over a range of spatial and temporal scales are of prime importance in  
620 assessing hydrosystem status. However, rivers are diverse and complex ecosystems, with  
621 significant variability of physical properties through both space and time. Future efforts  
622 involving detailed bathymetric data, light attenuation and water properties, may resolve depth-  
623 related confusion of SAV with substrate type, and are a prerequisite for multi-date vegetation  
624 monitoring, based on time series images.

625

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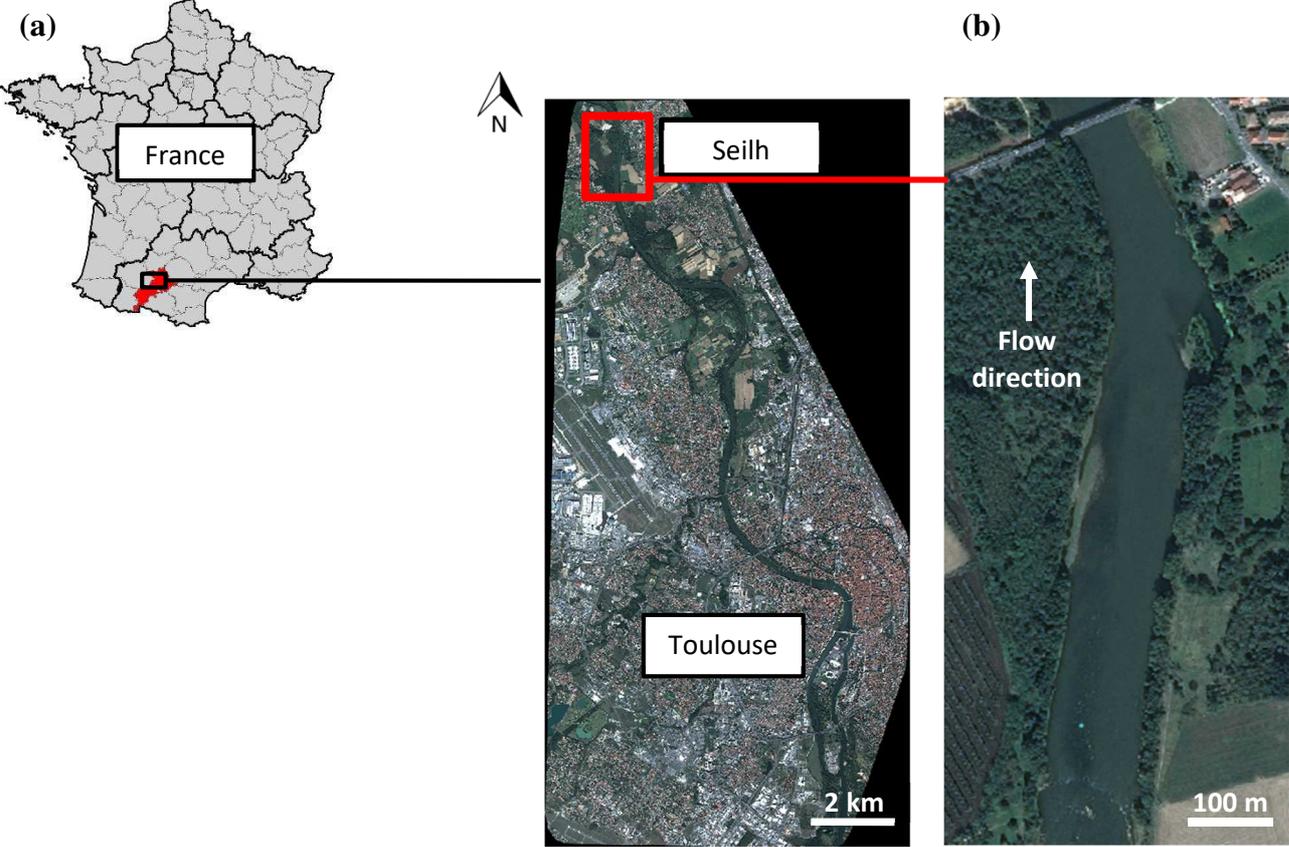
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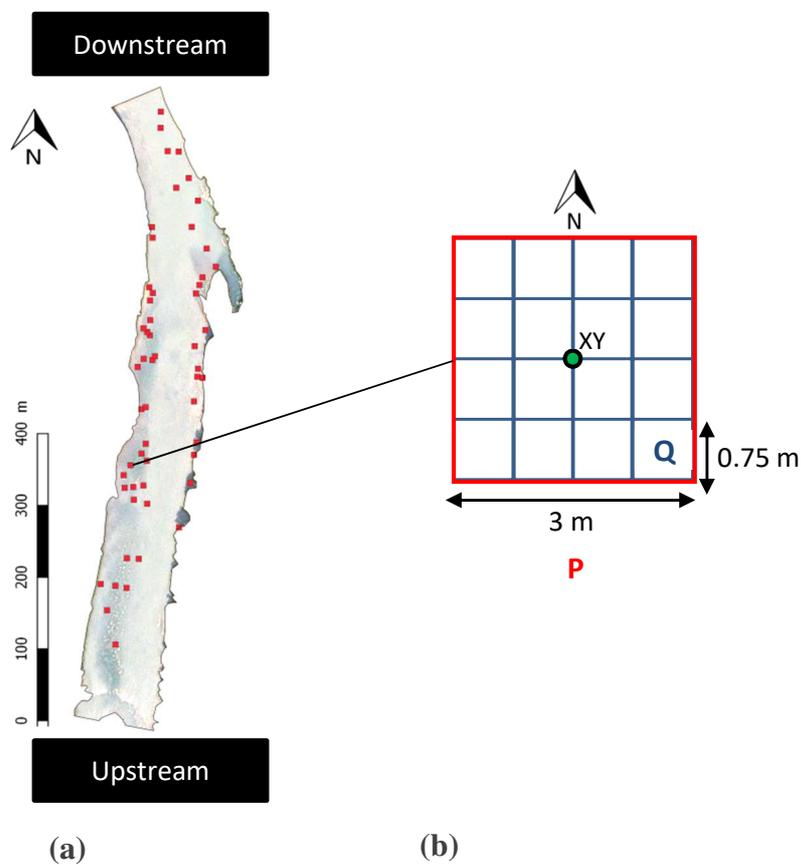
**Figure 1.** Study area. (a) Location of the study area in the southwest of France, north of Toulouse city on a raw Pléiades image; (b) Zoom in on a raw Pléiades image (50 cm of resolution) of the 1 km study site (Seilh).



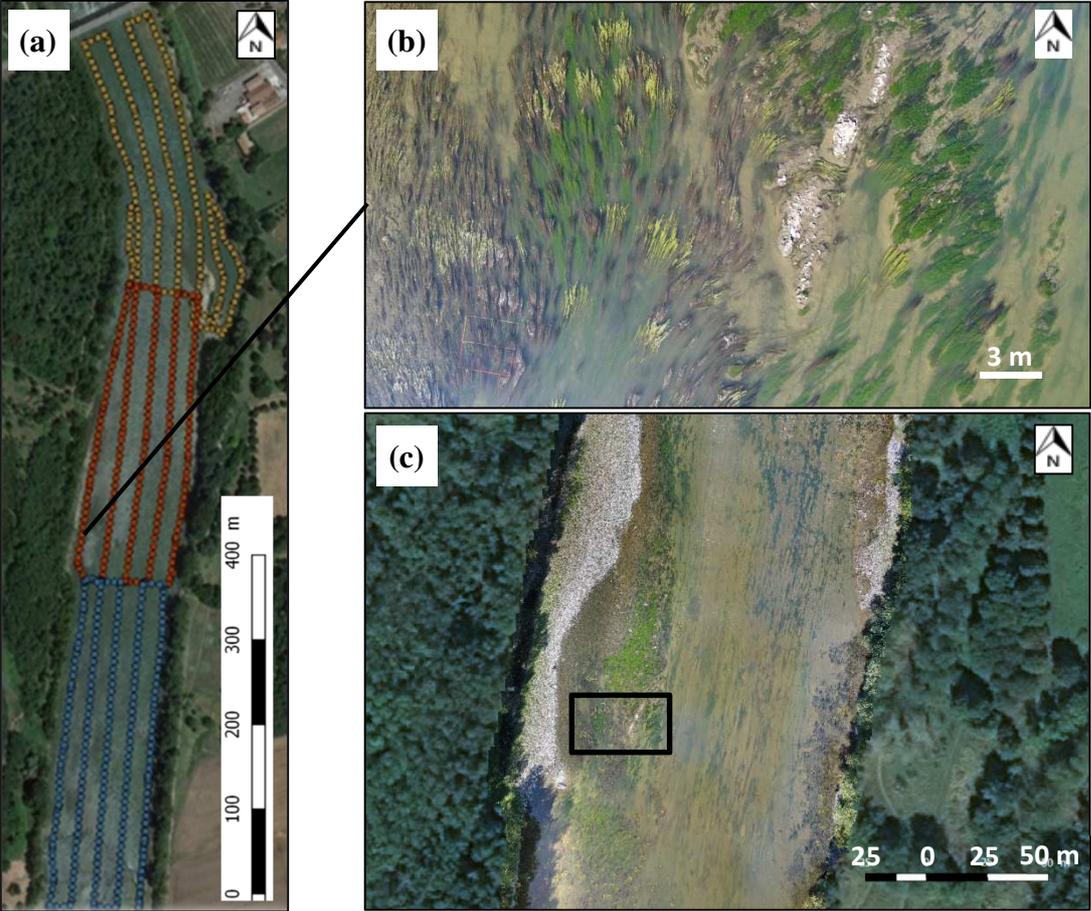
**Figure 2.** Sampling protocol for macrophyte monitoring.

(a) Location of the 55 sampling plots on the site (Seilh). The red squares indicate the geographical positions of the 55 field sampled plots (P).

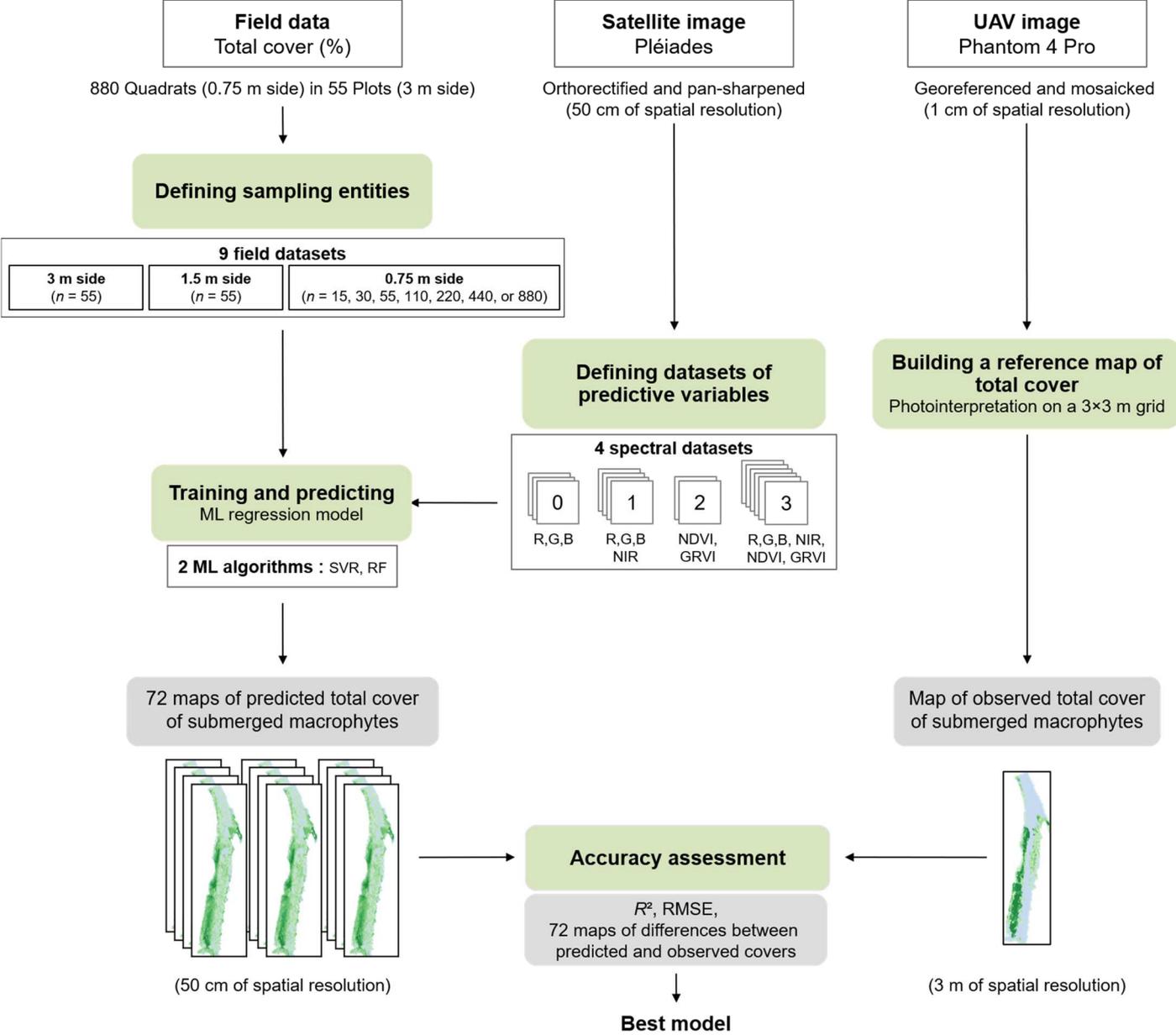
(b) Diagram of a plot (P): the red square represents the sampling plot of a 9 m<sup>2</sup> surface and the blue squares correspond to quadrats (Q). The green “XY” dot symbolizes where GPS coordinates were taken.



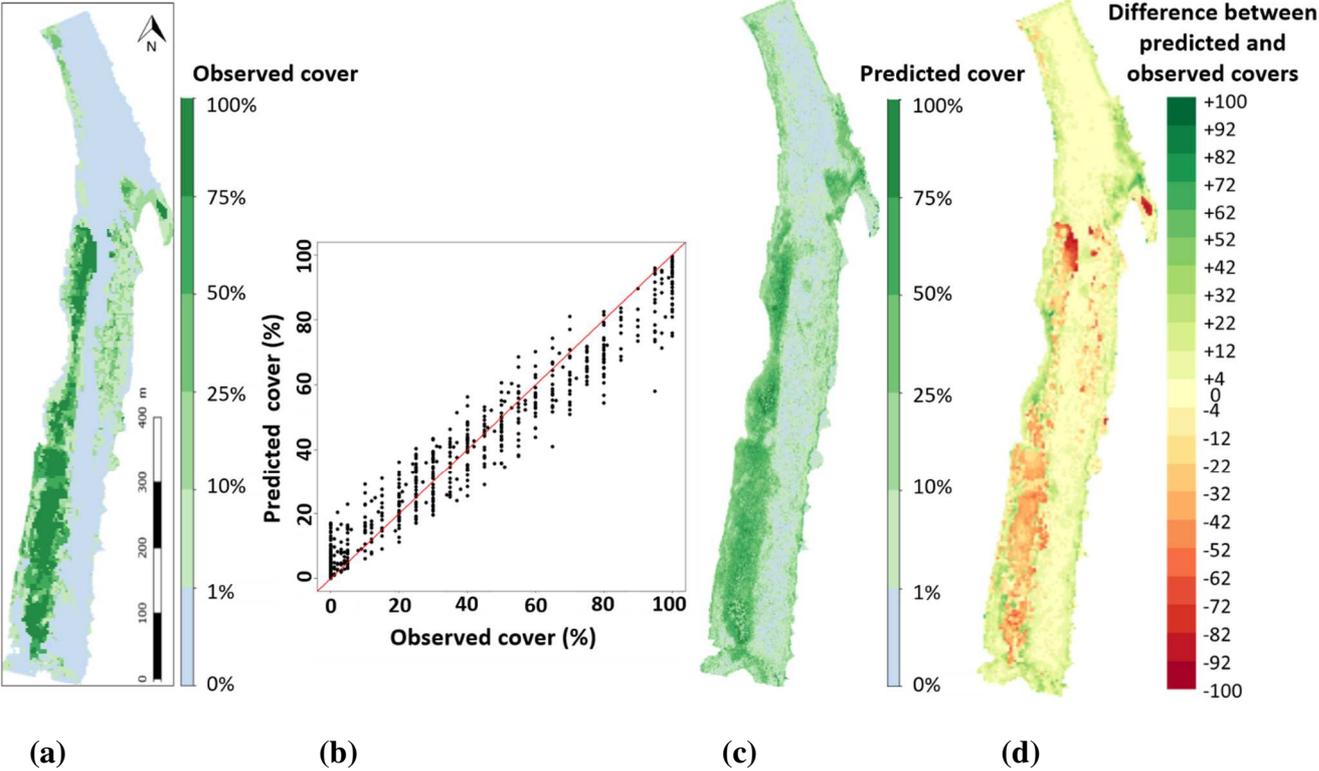
**Figure 3.** Drone images of the study site. (a) Image acquisition missions. Each point refers to a shooting, each colour corresponds to a mission; (b) Example of a raw drone image shot 30 m high above plurispecific macrophyte meadows; (c) Orthomosaic image fragment of the study site. The black rectangle represents the position of the individual shot in figure b.



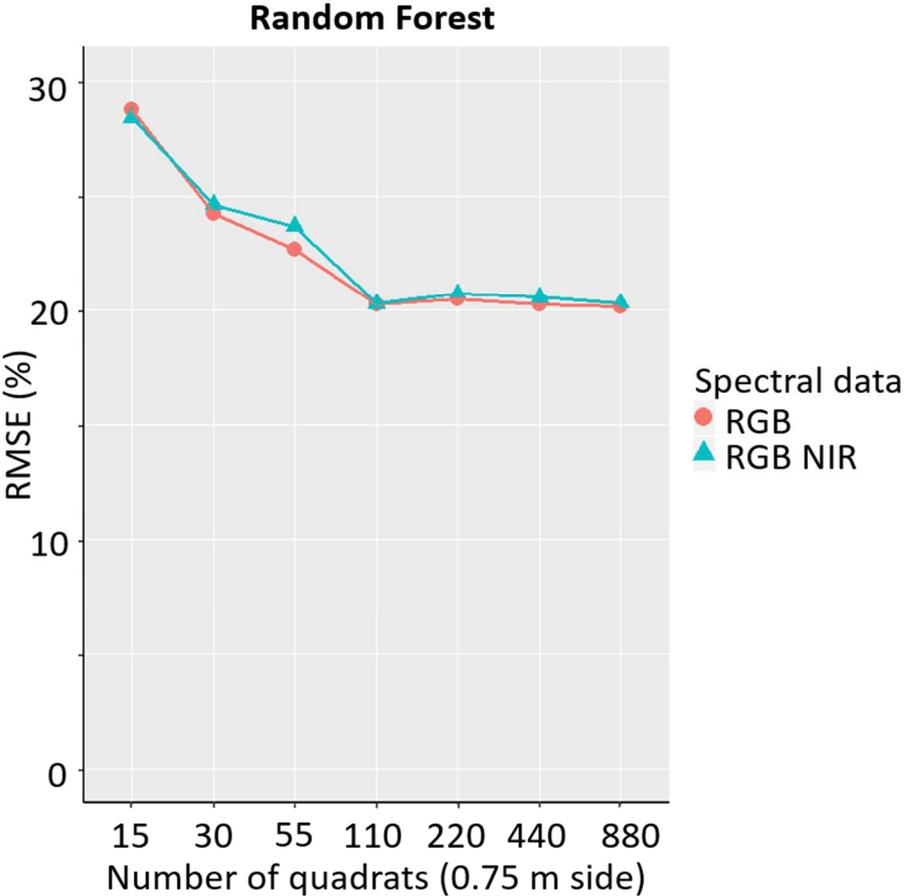
**Figure 4.** Flow chart illustrating the process to get total cover predictions from the Pléiades satellite image. Sampling entities can be 55 plots of 3 m side, 55 entities of 1.5 m side, or  $n$  quadrats of 0.75 m side (with  $n = 15, 30, 55, 110, 220, 440$  or  $880$ ).



**Figure 5.** Results from the Random Forest regression model using “dataset 1” (RGB and NIR) and based on 880 quadrats. (a) Observed cover map obtained from drone image photointerpretation (b) Predicted cover as a function of observed cover within quadrats; (c) Predicted cover map and (d) Difference map between predicted and observed covers.



**Figure 6.** Learning curves for the Random Forest regression model using datasets 0 (RGB) and 1 (RGB and NIR): RMSE as a function of the number of quadrats of 0.75 m side.



**Table 1.** Number of sampled entities according to macrophyte density ranges. Percentages in brackets represent the proportion of each cover group within the total number of samples.

<b>Macrophyte cover</b>	<b>Plots (P)</b>	<b>Quadrats (Q)</b>
0-1 %	16 (30 %)	359 (40 %)
1-10 %	6 (11 %)	65 (7 %)
11-25 %	7 (12 %)	99 (11 %)
26-50 %	12 (21 %)	160 (18 %)
51-75 %	10 (18 %)	100 (11 %)
>75 %	4 (7 %)	97 (11 %)
<b>Total number of entities</b>	<b>55</b>	<b>880</b>

**Table 2.** Prediction accuracy for different runs (differing by the algorithm, spectral dataset, or sampling level), assessed with the  $R^2$  coefficient of determination and the RMSE; only results involving datasets 0 and 1 are presented. This table is an extract of the most relevant results of table S2 (see Supporting Information).

Algorithm	Spectral data	Entity side size (m)	Number of entities	Maximum predicted cover (%)	$R^2$	RMSE (%)
<b>Random Forest</b>	RGB NIR (“dataset 1”)	3	55	87.3	0.69	21.28
		0.75	880	100.0	0.71	20.30
<b>Support Vector Regression</b>	RGB (“dataset 0”)	3	55	99.0	0.72	20.37
		0.75	880	100.0	0.74	19.40
	RGB NIR (“dataset 1”)	3	55	100.0	0.72	21.28
		0.75	880	100.0	0.67	21.29