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Fine-resolution landscape-scale biomass mapping using a spatiotemporal patchwork of LiDAR coverages

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ABSTRACT

Estimating forest aboveground biomass (AGB) at large scales and fine spatial resolutions has become increasingly important for greenhouse gas accounting, monitoring, and verification efforts to mitigate climate change. Airborne LiDAR is highly valuable for modeling attributes of forest structure including AGB, yet most LiDAR collections take place at local or regional scales covering irregular, non-contiguous footprints, resulting in a patchwork of different landscape segments at various points in time. Here, as part of a statewide forest carbon assessment for New York State (USA), we addressed common obstacles in leveraging a LiDAR patchwork for AGB mapping at landscape scales, including selection of training data, the investigation of regional or coverage specific patterns in prediction error, and map agreement with field inventory across multiple scales.

Three machine learning algorithms and an ensemble model were trained with Forest Inventory and Analysis (FIA) field measurements, airborne LiDAR, and topographic, climatic, and cadastral geodata. Using a novel set of plot selection criteria and growth adjustments to temporally align LiDAR coverages with FIA measurements, 801 FIA plots were selected with co-located point clouds drawn from a patchwork of 17 leaf-off LiDAR coverages (2014–2019). Our ensemble model was used to produce 30 m AGB prediction surfaces within a predictor-defined area of applicability (98% of LiDAR coverage) and within four vegetated landcover classes. The resulting AGB maps were compared with FIA plot-level and areal estimates at multiple scales of aggregation. Our model was overall accurate (% root mean squared error 22%–45%; mean absolute error 11.6–29.4 Mg ha⁻¹; mean error 2.4–6.3 Mg ha⁻¹), explained 73%–80% of field-observed variation, and yielded estimates that were largely consistent with FIA's design-based estimates (89% of estimates within FIA's 95% confidence interval). We share practical solutions to challenges faced in using spatiotemporal patchworks of LiDAR to meet growing needs for AGB prediction and mapping in support of broad-scale applications in forest carbon accounting and ecosystem stewardship.

1. Introduction

Mapping and monitoring forest aboveground biomass (AGB) has become increasingly important as the basis for large-scale accounting of carbon and greenhouse gas (GHG) fluxes in support of policy, regulatory, and land stewardship initiatives to mitigate global climate change. Although carbon stocks are the ultimate endpoint in GHG accounting, for several practical reasons forest AGB serves as a proxy for carbon stocks (and stock-changes) in large-scale accounting methodologies (Buendia et al., 2019; Woodall et al., 2015). Such applications, based primarily on field inventory, yield aggregate tabular estimates for political units (e.g., states, provinces), but offer little insights on landscape patterns within and across those units. Landscape-scale AGB maps with fine-resolutions can serve as inputs to GHG accounting and can help decision makers identify specific units of land for protection from deforestation, or as suitable candidates for reforestation, afforestation, or improved management in efforts to increase terrestrial carbon sequestration and offset GHG emissions from other sectors (Houghton et al., 2012; Houghton, 2005).

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National field sampling programs, like the United States Department of Agriculture's Forest Inventory and Analysis program (FIA) (Gray et al., 2012), provide estimates of AGB at individual plots and are scaled to larger areas through a design-based approach (Bechtold and Patterson, 2005). However, the resolution at which design-based estimates can be reliably produced is limited to relatively coarse units, such as states or counties, due to the density of the sample (McRoberts, 2011). For instance, the FIA program samples forest conditions and land cover across New York State (NYS) at a density of roughly one plot per 2,400 ha (Gray et al., 2012), while 65% of NY forestlands are owned in parcels smaller than 40 ha (L'Roe and Allred, 2013). Although FIA's extensive plot network was never designed to yield such localized results, this inherent resolution mismatch poses obstacles to accounting and decision-support applications that must account for local geography, both ecological and cadastral, to be practical and effective.

To address this need, model-based approaches combining field inventory data, like the FIA, with auxiliary remotely sensed data can produce predictions for all map units (pixels) in a given area. Airborne LiDAR has been established as a highly valuable remotely sensed data source for such purposes (Huang et al., 2019; Hurtt et al., 2019; Chen and McRoberts, 2016) offering detailed information on forest structure at fine spatial resolutions. However, LiDAR data are most commonly acquired at local to regional scales in irregular or noncontiguous footprints (Skowronski and Lister, 2012), resulting in a complex patchwork of data from component coverages acquired at different times with different sensors and mission parameters, in turn posing a host of challenges for broad-scale AGB mapping (Lu et al., 2014; Huang et al., 2019). These challenges include insufficient field inventory plots that spatially and temporally match LiDAR acquisitions and data discrepancies among LiDAR coverages.

Yet several groups have undertaken broad-scale AGB mapping efforts with LiDAR patchworks with varying degrees to which training data have been pooled from multiple LiDAR coverages. The choice between an individual or a pooled modeling approach often reflects practical considerations relating to sufficient sample size across all sub-regions and the cost of developing multiple models. Nilsson et al. (2017) did not pool at all, implementing a separate model trained for each coverage. Huang et al. (2019) pooled by ecoregion. Both Ayrey et al. (2021) and Hauglin et al. (2021) pooled all coverages but used a convolutional neural network and a mixed-effects model respectively, with differing protocols for inventory plot selection.

In this study, as part of a broader effort for map-based forest carbon accounting across NYS, we addressed several common challenges in using LiDAR patchworks for broad-scale, fine-resolution biomass modeling and mapping. We leveraged FIA inventories for model training and assessment data, and implemented FIA-developed methods to assess the agreement between our estimates and those produced by FIA. With the goal of producing a spatially explicit representation of FIA AGB information, we used a model-based approach to translate FIA's discrete plot-level estimates to wall-to-wall predictions at a 30 m resolution across a patchwork of 17 discrete LiDAR coverages in NYS.

We implemented a rigorous plot selection framework to limit temporal lags between LiDAR acquisitions and field inventories. When strict temporal alignment yielded too few plots, we leveraged repeated FIA inventories to boost the number of plots that temporally match LiDAR acquisitions without the incorporation of additional models or manual processes (Ayrey et al., 2021; Hauglin et al., 2021). In spite of this strategy, we were left with limited model training data, where some coverages and regions lacked sufficient information to support independent models, necessitating a single model or pooled approach.

We used machine learning (ML) algorithms including random forests (Breiman, 2001a), gradient boosting machines (Friedman, 2002), and support vector machines (Cortes and Vapnik, 1995), as well as 'stacked ensembles' of said algorithms (Wolpert, 1992). These have been shown to be better suited for pure prediction (Efron, 2020; Breiman, 2001b)

when compared to their conventional regression counterparts (Hauglin et al., 2021), especially when the input data are noisy and the processes governing the relationships between predictors and responses are complex or unknown, as is often the case in nature (Dormann et al., 2018). Furthermore, using pre-selected hyperparameters these algorithms can be trained in minutes with typical consumer-grade hardware, avoiding some of the drawbacks of computationally intense deep-learning approaches (Ayrey et al., 2021).

We also employed multiple strategies to address concerns that differences among sensors and mission parameters could lead to nonrandomly distributed errors in our AGB prediction surfaces. First, we produced an area of applicability (Meyer and Pebesma, 2021) mask to both examine the uniformity of our predictors across the component coverages, as well as to screen predictions based on anomalous predictor data. Second, we examined the spatial autocorrelation of our residuals and mapped our prediction error to identify the presence of region or coverage specific patterns. Finally, we assessed the agreement between our mapped predictions and FIA estimates across a range of scales (Riemann et al., 2010; Menlove and Healey, 2020).

2. Data and methods

2.1. LiDAR coverages

Our study relied upon a set of 17 LiDAR datasets hosted by the NYS GIS Program Office (GPO) covering 62.46% (7,835,773 ha) of NYS (hereafter "GPO-LiDAR area"; Table 1; Fig. 1). We selected individual coverages from the most recent five years of available data that contained temporally matching field data (2014–2019) to minimize sensor and data differences. All component coverages were originally collected to generate digital elevation models for flood risk analysis, and to this end were flown during leaf-off conditions. Several previous studies have shown that leaf-off LiDAR models can be as accurate as their leaf-on counterparts (Hawbaker et al., 2010; White et al., 2015; Anderson and Bolstad, 2013). Further specifications for each LiDAR coverage can be found in Supplementary Materials S1.

2.2. Field data

Two field datasets were compiled from the USDA FIA inventory in NYS with the distinct purposes of model development and map assessment. The FIA program compiled AGB estimates for trees \geq 12.7 cm (5 in) diameter at breast height (Gray et al., 2012), and were converted to units of megagrams per hectare (Mg ha⁻¹). The FIA uses permanent inventory plots arranged in a quasi-systematic hexagonal grid that are remeasured on a rolling 5-7 year basis in the Northeastern United States (Bechtold and Patterson, 2005). Tree measurements, and subsequently AGB estimates, are only recorded on portions of plots considered forested. For an area to be considered forested by the FIA, the area must be at least 10% stocked with trees, at least 0.4 ha (1 acre) in size, and at least 36.58 m (120 ft) wide. Additionally, any lands meeting these minimum requirements, but developed for nonforest land uses, are not considered forested. With an understanding that some nonforest conditions contained AGB that was not measured by FIA, we assumed that any nonforest condition contained 0 AGB.

FIA plots are composed of four identical circular subplots with radii of 7.32 m (24 ft), with one subplot centered at the macroplot centroid and three subplots located 36.6 m (120 ft) away at azimuths of 360°, 120°, and 240° (Bechtold and Patterson, 2005). The plot locations were provided by the FIA program in the form of average coordinates, collected over multiple repeat visits, representing the centroid of the center subplot, which we then used to build a polygon dataset representing the entire plot layout including all four subplots (Fig. 2). Averaged coordinates were necessary due to the lack of precision of initial GPS coordinates for the macroplot centroids (Hoppus and Lister, 2005). Any reference to an FIA plot hereafter implies the aggregation of all four component subplots.

Component LiDAR coverage metadata. IDs for cross figure correspondence; Year of acquisition; Area covered (ha); Pulse density (PD) in pulses per m^2 ; Number of FIA plots in model and map assessment datasets. Area of Applicability (AOA) in percent of LiDAR coverage pixels considered inside of the area of applicability. AOA computation conducted after initial Land Change Monitoring, Assessment, and Projection (LCMAP) primary classification masking.

ID	Coverage name	Year	Area	PD	% AOA	Model	Assessment
1	Erie, Genesee & Livingston	2019	555,853	2.04	97.72	11	33
2	Fulton, Saratoga, Herkimer & Franklin	2018	557,421	2.60	99.61	39	84
3	Southwest B	2018	527,075	1.98	98.06	28	66
4	Cayuga & Oswego	2018	438,201	2.78	96.89	25	43
5	Southwest	2017	423,714	1.98	98.53	33	74
6	Franklin & St. Lawrence	2017	977,620	2.69	99.24	104	188
7	Oneida Subbasin	2017	264,886	2.10	96.02	15	31
8	Allegany & Steuben	2016	309,081	1.69	97.77	32	53
9	Columbia & Rensselaer	2016	248,839	1.69	98.19	17	28
10	Clinton, Essex & Franklin	2015	600,755	2.23	98.70	115	127
11	Warren, Washington & Essex	2015	611,704	3.24	99.37	106	128
12	Madison & Otsego	2015	471,564	2.13	99.24	56	92
13	3 County	2014	755,629	2.04	96.12	92	114
14	Long Island	2014	315,542	2.04	94.77	22	12
15	Schoharie	2014	256,464	2.04	95.12	31	40
16	New York City (NYC)	2014	77,211	1.54	90.40	2	1
17	Great Lakes	2014	444,215	2.04	98.47	73	103
	GPO-LiDAR		7,835,773		98.12	801	1217

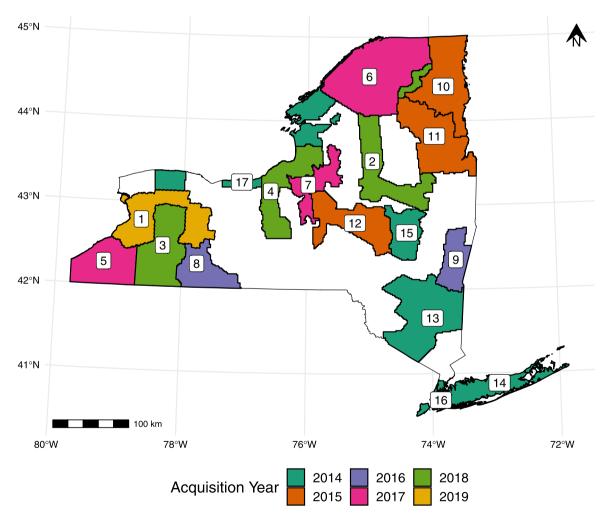


Fig. 1. GPO-LiDAR component coverages, colored by year of acquisition and labeled by ID numbers in Table 1.

2.2.1. Model development dataset

In selecting plots for model development (model dataset hereafter; Table 1) we aimed to maximize the number of reference plots (Fassnacht et al., 2014), while minimizing the temporal lag between LiDAR acquisitions and inventories and ensuring high quality co-registered LiDAR data (White et al., 2013; CEOS, 2021). Temporal misalignment between plots and LiDAR has been shown to introduce error in model predictions; Gonçalves et al. (2017) found that 7%–17% of their prediction error could be attributed to 3-year lags between inventories and LiDAR acquisitions.

Summary of plot selection criteria for the model dataset.

	Criteria	Description	Num plots
	1	Plots inventoried in the same	324
Include		year as a LiDAR acquisition.	
	2	Plots growth-adjusted to	843
		temporally align with LiDAR	
		acquisition.	
	3	Plots where measured $AGB = 0$	327
		Mg ha ⁻¹ but maximum LiDAR	
		return > 1 m.	
	4	Plots where the convex hull of	13
Exclude		LiDAR returns clipped to a given	
		subplot did not contain at least	
		90% of the subplot area.	
	5	Plots duplicated in neighboring	26
		coverages. Temporal matches	
		given preference, followed by	
		inventory recency.	

We only considered FIA plots when all subplots were marked as measured and when forest conditions were uniform at the plot level (all subplots completely forested or completely nonforested). Only 324 plots were available when we required a strict temporal match between inventories and LiDAR acquisitions (Table 2; Criterion 1). To increase the number of reference plots, we applied a growth adjustment to FIA plots with inventories both before and after the LiDAR acquisition, or 'bracketing inventories'. For plots with bracketing inventories, we excluded those where AGB decreased by \geq 5% between measurements indicating a disturbance event on the plot. For the remaining (relatively undisturbed) plots with bracketing inventories, AGB at the time of the LiDAR collection was then estimated by linearly interpolating between bracketing FIA estimates. This procedure added 843 plots to our dataset (Table 2; Criterion 2), and is analogous to existing growth adjustment methods (Gonçalves et al., 2017; Gobakken and Næsset, 2008).

Several plot exclusion rules were implemented to filter duplicates and remove problematic observations resulting from vegetation in nonforest conditions, interfering structures, or other data anomalies which would otherwise degrade the relationships between AGB and our predictor variables (Table 2; Criteria 3–5). Point clouds were clipped to the constructed plot polygons and were excluded with criteria 3–5 (Fig. 2).

2.2.2. Map assessment dataset

For our map assessment dataset (assessment dataset hereafter; Table 1) we were primarily concerned with maintaining FIA's probability sample of inventory plots in order to leverage unbiased estimators of map agreement metrics (Stehman and Foody, 2019; Riemann et al., 2010). As in the model dataset, only those plots with all four subplots marked as measured were considered. Excluding non-measured plots, however, does not invalidate FIA's probability sample, as the FIA program assumes these plots to be randomly distributed across the landscape (Bechtold and Patterson, 2005). To this end, we selected FIA plots on a coverage-by-coverage basis (Fig. 1), choosing only plots that had been inventoried in +/-2 years from the year of the LiDAR acquisition, but not inventoried in the same year as the LiDAR acquisition to maintain independence from the model dataset (Table 3; Criterion 1). Where plots were shared with the model dataset, AGB values differed, with AGB values derived from measurements in the map assessment dataset, and AGB values derived from our growth-adjustment process in the model dataset, therefore representing different points in time. Furthermore, plots falling outside of our mapped area, based on our landcover and area-of-applicability masks (Section 2.6), were excluded as they were considered outside our population of interest (Table 3; Criteria 2-3).

Table 3

Summary of plot selection criteria for the assessment dataset.

	Criteria	Description	Num plots
Include	1	Plots inventoried in $+/-2$ years of LiDAR acquisition but not in the same year as the acquisition.	1639
Exclude	2	Plots with any portion of the plot footprint outside of our landcover mask.	291
	3	Plots with any portion of the plot footprint outside of our area of applicability mask.	131

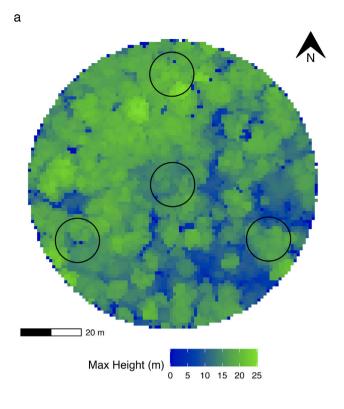
2.3. LiDAR and auxiliary data pre-processing

First, we height-normalized the raw LiDAR data using digital terrain models (DTM) computed from each LiDAR coverage separately, where each DTM was computed on the fly, such that an elevation value was imputed for each point in the raw LiDAR point cloud, and a complete DTM surface was never created or saved. The point-wise elevations were estimated using an inverse-distance-weighted imputation using pre-classified ground and water returns with five neighbors and a power of two. From the height-normalized data we computed 40 predictors (Supplementary Materials S2) based on previous studies (Hawbaker et al., 2010; Huang et al., 2019; Pflugmacher et al., 2014). Predictors at individual FIA plots were computed from LiDAR returns clipped to only the measured subplot areas and then pooled at the plot level (Fig. 2). The corresponding predictors computed for map pixels were based on the set of returns within each 30 m cell. The lidR (Roussel and Auty, 2020; Roussel et al., 2020) package in R (R Core Team, 2021) was used for height-normalization and predictor generation.

A group of steady-state predictors was included to represent geospatial variation in climate and topography (Kennedy et al., 2018). Additionally, a 2019 tax parcel layer was incorporated as a set of boolean indicator variables (Supplementary Materials S3). Tax codes and categories provide cadastral information related to land-use and management (Thompson et al., 2011). All auxiliary predictors were reprojected and resampled to pixel geometries matching the 30 m LiDAR predictor surfaces.

2.4. Model development

Three ML models were fit to a randomly selected 80% of the model dataset (training set; n = 630), leaving the 20% remaining plots as an independent testing set to assess the final model performance on point clouds clipped to FIA plot polygons (testing set; n = 171). All predictors described in Section 2.3 were included in the model



b

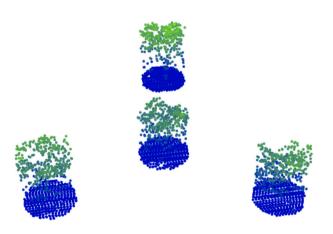


Fig. 2. (a) FIA plot layout overlaid on a 1 m LiDAR-derived max-height surface. (b) Height-normalized LiDAR returns clipped to individual FIA subplots and pooled prior to predictor computation.

training process without applying any dimensionality reduction procedures, and the following ML models with respective R packages were used: Random forest (RF), ranger package (Wright and Ziegler, 2017), stochastic gradient boosting machines (GBM), lightgbm package (Ke et al., 2021, 2017), and support vector machines (SVM), kernlab package (Karatzoglou et al., 2004). Each of the models were tuned using the 80% training dataset and an iterative grid search, starting by testing wide ranges of hyperparameters using five-fold cross validation and then narrowing down to only the most performant combinations over several iterations. Models then used the most accurate sets of hyperparameters in all other analyses. For each of the *n* observation in the training dataset, all three component models were fit, using their optimal hyperparameters, with n - 1 observations. Predictions for each component model were made for the *n*th (left out) observation. A linear regression model was used to estimate AGB as a function of these leave-one-out predictions, combining the three ML models in a stacked ensemble to better reflect model selection uncertainty (Wintle et al., 2003; Dormann et al., 2018), and to reduce the generalization error of our component models (Wolpert, 1992). The linear regression ensemble was constructed as follows:

$$AGB = \beta_0 + \beta_1 \cdot RF + \beta_2 \cdot LGB + \beta_3 \cdot SVM + \beta_4 \cdot RF \cdot LGB + \beta_5 \cdot RF \cdot SVM + \beta_6 \cdot LGB \cdot SVM + \beta_7 \cdot RF \cdot LGB \cdot SVM$$
(1)

where β_* are coefficients estimated through ordinary least squares regression, and *RF*, *GBM*, and *SVM* are the respective component model predictions. The selected hyperparameters for each component model and the coefficients in the linear regression ensemble are available in Supplementary Materials S4.

2.5. Model performance

Model performance was assessed against the 20% testing partition of the model dataset (Table 1) based on metrics including root-meansquared error in Mg ha⁻¹ (RMSE, Eq. (2)), percent RMSE (% RMSE, Eq. (3)), mean absolute error in Mg ha⁻¹ (MAE, Eq. (4)), percent MAE (% MAE, Eq. (5)), mean error in Mg ha⁻¹ (ME, Eq. (6)), and the coefficient of determination (R^2 , Eq. (7)) as follows:

RMSE =
$$\sqrt{(\frac{1}{n})\sum_{i=1}^{n}(y_i - \hat{y}_i)^2}$$
 (2)

$$\% \text{ RMSE} = 100 \cdot \frac{\text{RMSE}}{\bar{y}}$$
(3)

MAE =
$$(\frac{1}{n}) \sum_{i=1}^{n} (|y_i - \hat{y}_i|)$$
 (4)

$$\% \text{ MAE} = 100 \cdot \frac{\text{MAE}}{\bar{y}} \tag{5}$$

$$ME = (\frac{1}{n}) \sum_{i=1}^{n} (y_i - \hat{y}_i)$$
(6)

$$R^{2} = 1 - \frac{\sum_{i=1}^{n} (y_{i} - \hat{y}_{i})^{2}}{\sum_{i=1}^{n} (y_{i} - \bar{y})^{2}}$$
(7)

where *n* is the number of FIA plots included in the data set, \hat{y}_i is the predicted value of AGB, y_i the AGB value measured at the corresponding location, and \bar{y} the mean AGB value from FIA field measurements.

Standard errors for R^2 and RMSE were computed as follows:

$$SE_{boot} = \sqrt{\frac{Var_{boot}}{n}}$$
(8)

where *n* is the number of FIA plots included in the dataset, and Var_{boot} is computed as the variance of R² and RMSE estimates for 1000 iterations of bootstrap resampling. Standard errors for MAE and ME were computed as follows:

SE =
$$\sqrt{\frac{\sum_{i=1}^{n} (e_i - \bar{e})^2}{n-1}}$$
 (9)

where e_i is the error at an FIA plot, and \bar{e} is the mean error from all FIA plots included. In the case of MAE, $e_i = |y_i - \hat{y_i}|$, the absolute value of the error at a given FIA plot.

2.6. AGB mapping and post-processing

The linear model ensemble was used to make predictions for all 30 m pixels within the GPO-LiDAR area. Predictions on newer coverages superseded those based on older coverages in areas where neighboring LiDAR coverages overlapped. All AGB prediction surfaces were projected to match Landsat 30 m pixel geometries to avoid mixed pixel effects in subsequent raster overlay analyses (Wulder et al., 2022).

With recognition that our predictions are best suited to areas populated by woody biomass, we tabulated our predictions across landcover types determined from the United States Geological Survey's Land Change Monitoring, Assessment, and Projection (LCMAP) primary classification products (Brown et al., 2020; Zhu and Woodcock, 2014), which has a reported overall accuracy of 77.4% in the eastern United States for the years 1985–2018 (Pengra et al., 2020). LCMAP's annual resolution (1985–2019) allowed for temporal alignment with the patchwork of LiDAR-AGB surfaces, and shared identical pixel geometries. We masked our AGB prediction surfaces to remove Developed, Water, and Barren pixels, and then tabulated AGB by the four remaining vegetated LCMAP classes of Tree cover, Grass/Shrub, Cropland, and Wetland.

Lastly, we computed an area of applicability (AOA) surface as a final mask for our AGB prediction surfaces (Meyer and Pebesma, 2021; Supplementary Materials S5). This algorithm calculates distance in predictor space, using the full set of predictors available to our component ML models (Section 2.3), to identify pixels containing predictor data that are not sufficiently well-represented in the training data. We used this mask to remove pixels falling outside of our AOA, and thus restricted our maps to only areas where we can expect our models to perform similarly to training and test set performance.

2.7. Map agreement assessment

We assessed the agreement between our AGB maps and FIA reference data following approaches prescribed by Riemann et al. (2010) and Menlove and Healey (2020). The former evaluated agreement across a range of scales and accounts for the mismatch in spatial support between map aggregate estimates (many pixels) and FIA aggregate estimates (few plots) by only extracting pixels coincident with FIA plots. The latter compared FIA-derived AGB estimates — which have been adjusted for forest cover within, and area-extrapolated to, hexagon map units — to zonal averages of our mapped AGB.

Following Riemann et al. (2010) we compared our AGB prediction surfaces to the assessment dataset (Section 2.2). Comparisons were made at both the plot-to-pixel scale and within variably-sized hexagons with distances between centroids ranging from 10 km (8660 ha) to 100 km (866.025 ha). As an extension of the Riemann et al. (2010) methodology we assessed prediction error (RMSE, MAE, ME) with choropleth maps that summarized the mapped residuals and FIA reference data distributions within hexagon units with centroids spaced 50 km apart. We also grouped plot to pixel results by the majority LCMAP classification at each plot, to demonstrate the level of agreement across vegetated landcover classes. To evaluate the presence of spatial autocorrelation among mapped residuals, which could indicate the presence of regional or coverage specific error in our model's predictions, Global Moran's I statistics were computed for search radii ranging from 1 to 50 km using both the model dataset and the assessment dataset separately (Moran, 1950).

Following the Menlove and Healey (2020) approach we compared the average of our masked predictions, weighted by the proportion of each pixel intersecting a given hexagon, to a set of FIA-derived estimates for 64,000 ha hexagons representing FIA's finest acceptable scale for the most recent inventory cycle in NYS (2013–2019) (Menlove and Healey, 2020). As recommended by Menlove and Healey (2020), we accounted for differences in forest definitions between the FIA estimates and our mapped estimates by dividing FIA estimates by the total area of vegetated (based on LCMAP Tree cover, Wetland, Cropland, Grass/Shrub) pixels within each hexagon. Lastly, we limited this comparison to only those hexagons that contained > 10% mapped area, using the summed area of all pixels inside of our LCMAP and AOA mask (Section 2.6) to define mapped area within each hexagon.

The exactextractr (Baston, 2021), spdep (Bivand et al., 2013), sf (Pebesma, 2018), raster (Hijmans, 2021), and terra (Hijmans, 2022) packages in R (R Core Team, 2021) were used to conduct all analyses described here. Further description of this assessment is included in Supplementary Materials S6.

Table 4

Model performance metrics (as defined in Section 2.5) against 20% testing partition of
the model dataset (n = 171). RMSE, MAE, and ME in Mgha ⁻¹ .

	RF	GBM	SVM	Ensemble
RMSE	40.55	41.02	41.63	41.07
% RMSE	37.34	37.78	38.34	37.82
MAE	29.08	29.68	29.27	29.36
ME	3.14	1.05	-1.90	3.10
R ²	0.75	0.74	0.73	0.74

3. Results

3.1. Growth-adjusted field plots

Of the 801 FIA plots in our model dataset, 562 were growthadjusted, representing about 70% of the full dataset. The average annual AGB increment applied to all growth-adjusted plots was 1.84 Mg ha⁻¹ year⁻¹ over an average of 3.96 years. The largest annual AGB increment applied was 11.17 Mg ha⁻¹ year⁻¹ over 3 years. The longest adjustments (applied to 5 plots) occurred over 10 years at an average annual increment of 2.63 Mg ha⁻¹ year⁻¹.

3.2. Model performance

The three component ML models and the ensemble model were generally accurate against the 20% test partition of the model dataset (Table 4), but were more accurate towards the mean of the reference distribution (Fig. 3). This resulted in slight overpredictions at smaller AGB plots and underpredictions at larger AGB plots (Fig. 3). While there were only marginal differences between the model performance metrics, and RF and GBM yielded slightly smaller RMSE values than the ensemble model (Table 4), we selected the ensemble model for mapping and all further analysis based on first principles described in Section 2.4, and its improved ability to make predictions towards the extremes of the reference distribution relative to any single component model.

3.3. AGB by landcover class

Approximately 62% of all pixels, accounting for 87% of the total mapped AGB (Fig. 4), were contained within areas identified as Tree cover by Land Change Monitoring, Assessment, and Projection (LCMAP) primary classification products (Table 5). All other LCMAP classes contained comparatively small, but non-zero estimates of AGB, with Wetlands and Grass/Shrub containing larger average predictions within small portions of the mapped area, and Cropland containing smaller average predictions across a larger proportion of the mapped area (Table 5).

3.4. Area of applicability

All mapped (vegetated) LCMAP classes contained AOA \geq 97.64% and the AOA breakdown across component LiDAR coverages (after LCMAP masking) was uniform (\geq 90.4%) with NYC and Long Island the only two coverages with \leq 95% AOA (Tables 1 and 5). In total, 98.12% of the GPO-LiDAR area was considered inside the AOA after initial LCMAP masking.

3.5. Map agreement assessment

We observed improved agreement between the mapped estimates and the FIA estimates as the aggregation unit size increased, with %

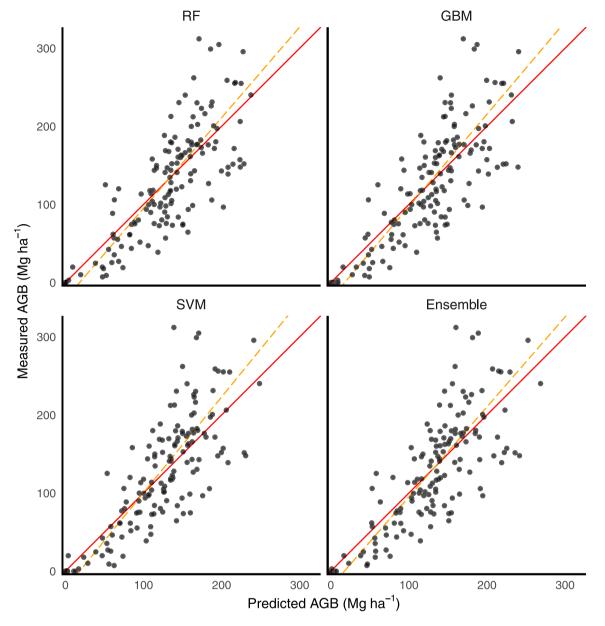


Fig. 3. Measured vs predicted AGB scatter plot for the 20% testing portion of the model dataset. AGB values in Mg ha⁻¹. Geometric mean functional relationship (GMFR) trend line shown with dashed (orange) line, and 1:1 line shown with solid (red) line.

Summary of FIA reference plots in the model dataset and mapped predictions by LCMAP landcover classes. n = number of plots. Mean AGB values in Mg ha⁻¹. Total AGB values in millions of metric tons. Area in hectares. AOA expressed as percent of LCMAP classified pixels considered inside of the area of applicability.

LCMAP	Reference plots		ference plots Mapped					
	n	Mean AGB	Area	% Area	Mean AGB	Total AGB	% AGB	% AOA
Tree cover	599	137.97	4,251,812	62.42	132.66	564.06	87.38	98.13
Cropland	132	1.98	1,748,905	25.68	14.11	24.67	3.82	98.25
Wetland	36	115.13	613,578	9.01	77.22	47.38	7.34	97.64
Grass/Shrub	12	30.68	196,852	2.89	47.86	9.42	1.46	98.17

RMSE decreasing from 45 to 15% and MAE decreasing from 28 to 10 Mg ha⁻¹ (Table 6, Figs. 5 and 6). The scatter plot results (Fig. 5) for the plot to pixel and 10 km scales show a pattern of large overpredictions where FIA reference values are 0 Mg ha⁻¹ AGB — likely an indication of the structural zeroes introduced by FIA's strict definition of forest. Notably, ME increased as a function of aggregation unit size, which appears to be a reflection of outlier hexagons with very small samples

rather than any additive effects from aggregation (Supplementary Materials S6). We again observed less accurate predictions at the extremes of the distributions across most scales, though our extreme predictions (small and large) became more accurate as a function of aggregation unit size, as indicated by the geometric mean functional relationship (GMFR; Riemann et al., 2010) slope approaching 1 at the largest scales (Figs. 5 and 6).

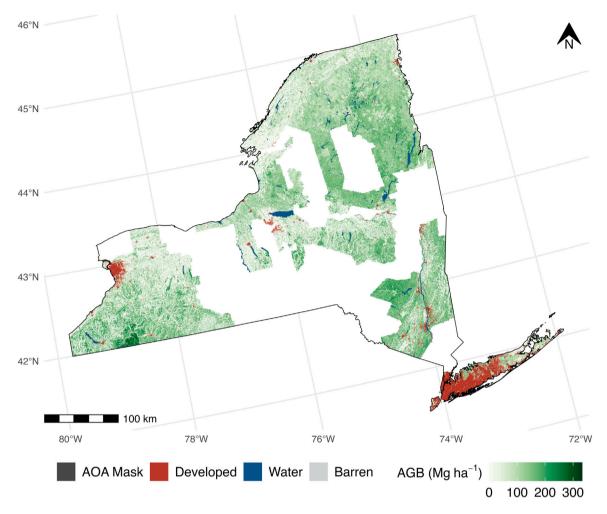


Fig. 4. AGB prediction surfaces reflecting a temporal patchwork of conditions across the 17 component LiDAR coverages in the GPO-LiDAR area with New York State (USA).

The model's mapped residuals grouped to 50 km spaced hexagons did not reveal any observable spatial patterns of prediction error (RMSE, MAE, ME) (Fig. 7). The ME map (Fig. 7c) was dominated by near-zero positive prediction error but indicated the tendency for hexagons with negative prediction ME to have larger mean FIA AGB values, while hexagons with positive ME were more likely to have smaller mean FIA AGB values (Fig. 7g). The RMSE map (Fig. 7a) and MAE map (Fig. 7b) largely mirrored one another.

Comparison of our AGB maps with FIA's design based estimates of AGB density (Menlove and Healey, 2020) indicated overall strong agreement, with 89% of our estimates falling within the FIA estimate 95% confidence intervals (CI) (Fig. 8). The majority of our estimates falling outside the CI were at the lower range of the AGB distribution, reinforcing the observed pattern of overprediction in comparison to FIA-derived estimates at the low end of the reference distribution.

Map agreement across LCMAP vegetated classes was highly variable, with the smallest % RMSE for plots classified as Tree cover by LCMAP (Table 7). Absolute RMSE for Grass/Shrub and Cropland plots was small (< 24 Mg ha⁻¹), but their relative values were quite large (% RMSE > 100%), reflecting the small FIA AGB averages within these classes. Mean errors were positive and large for Grass/Shrub and Croplands (14.64 Mg ha⁻¹ and 6.67 Mg ha⁻¹ respectively; Table 7), suggesting that errors contained in these two landcover classes had a relatively large contribution to the overall patterns of positive ME and overprediction on the low end of the FIA AGB distribution across the entire GPO-LiDAR area.

The global Moran's I analysis with the assessment dataset found evidence of very weak spatial autocorrelation (≤ 0.10) in mapped

residuals for all search radii (Supplementary Materials S7). Notably, when the analysis was conducted with the model dataset, similarly very weak spatial autocorrelation (≤ 0.08) was only evident for search radii of 9 km to 14 km, but was not evident at any other scales (Supplementary Materials S7).

4. Discussion

In this study we attempted to use a patchwork of 17 discrete LiDAR coverages for broad-scale, fine-resolution forest aboveground biomass (AGB) mapping across New York State (NYS), USA. Faced with a limited sample of temporally aligned field inventory data, we leveraged repeated inventories to boost the model training sample, and used a machine learning ensemble model to produce accurate predictions. We addressed concerns of sensor and mission discrepancies among component LiDAR coverages by investigating spatial patterns of prediction error, and by using an AOA mask to both show predictor uniformity across coverages, as well as to mask predictions based on anomalous data. Our results demonstrated that our maps accurately characterized the spatial patterns of AGB across the state, and showed that our predictions have strong agreement with FIA estimates across a range of aggregation scales.

4.1. Growth-adjusted field plots

Our approach to solve the common lack of temporally coincident LiDAR and field data was parsimonious and effectively tripled the sample size while achieving accurate modeling results. We did this by

Map agreement results for select scales. Distance = distance between hexagon centroids in km; PPH = plots per hexagon; n = number of comparison units (plots or hexagons); RMSE, MAE, ME in Mg ha⁻¹. All accuracy metrics as defined in Section 2.5. Standard errors in parentheses.

Distance	n	PPH	% RMSE	RMSE	MAE	ME	R ²
plot:pixel	1217	-	44.87	40.93 (0.04)	28.12 (0.85)	4.40 (1.17)	0.73 (0.01)
10	739	1.65	37.95	34.06 (0.05)	23.95 (0.89)	4.63 (1.24)	0.77 (0.01)
25	199	6.12	27.95	24.68 (0.13)	17.13 (1.26)	2.36 (1.75)	0.80 (0.01)
50	72	16.90	25.86	21.17 (0.41)	14.26 (1.86)	6.26 (2.40)	0.78 (0.01)
100	26	46.81	22.28	17.39 (0.71)	11.58 (2.59)	5.42 (3.30)	0.74 (0.03)

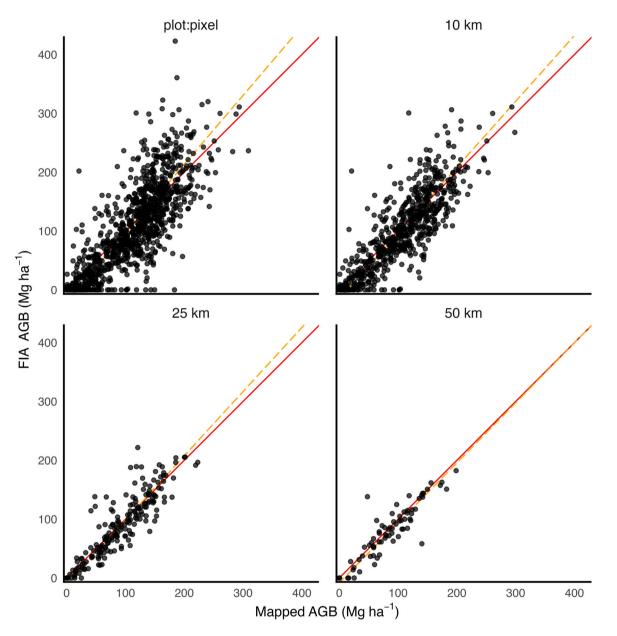


Fig. 5. Comparing mapped AGB to FIA estimated AGB across selected scales represented by distances between hexagon centroids (plot:pixel, 10 km, 25 km, and 50 km). AGB values in Mg ha⁻¹. GMFR trend line shown with dashed (orange) line, and 1:1 line shown with solid (red) line.

Map agreement at the plot to pixel scale, grouped by LCMAP classification. n = number of plots; RMSE, MAE, ME in Mg ha⁻¹. All accuracy metrics as defined in Section 2.5. Standard errors in parentheses, with a minimum of 0.01 for display.

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LCMAP	n	% RMSE	RMSE	MAE	ME	R ²
Tree cover	797	36.35	45.85 (2.56)	34.42 (1.07)	3.83 (1.62)	0.55 (0.01)
Cropland	303	229.82	18.90 (2.40)	10.32 (0.91)	6.67 (1.02)	0.40 (0.03)
Wetland	91	62.48	51.36 (30.77)	35.56 (3.91)	-1.13 (5.41)	0.55 (0.01)
Grass/Shrub	26	119.65	23.71 (17.01)	16.03 (3.49)	14.64 (3.73)	0.50 (1.24)

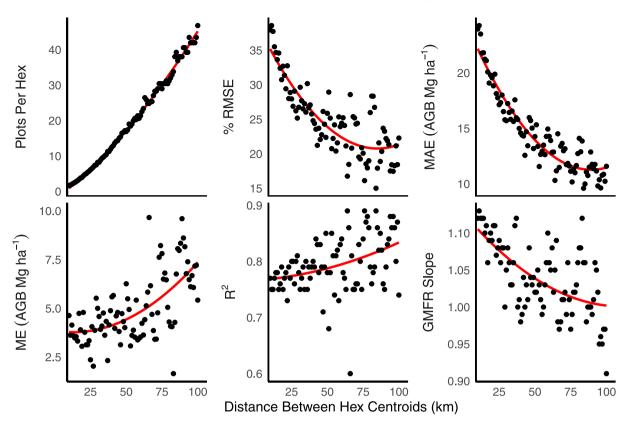


Fig. 6. Summary assessment metrics (as defined in Section 2.5) comparing mapped predictions to FIA estimates as a function of aggregation unit size (described by distances between hexagon centroids). Red trend lines produced using quadratic regression.

leveraging the existing inventory data, without additional field campaigns, remotely sensed data, or growth and yield models. However, a potential limitation to this approach is the requirement of regular historical inventories so that bracketing inventory years can be identified for LiDAR acquisitions. Additionally, the maximum temporal distance between growth-adjusted AGB values and the nearest measured AGB values in our dataset was 10 years; growth-adjustment following our approach may not be reliable for longer temporal lags.

4.2. Pooled modeling

Despite our efforts to boost the amount of training information available to models via growth adjustment, we were left with a nonuniform spatial arrangement of FIA plots across the GPO-LiDAR area. Several coverages contained fewer than 20 plots in the model dataset (Table 1), requiring a pooled modeling approach. Pooling information from all component LiDAR coverages allowed our model to borrow information from coverages with more training data to build relationships between our predictors and AGB that were then applied to coverages with less training data.

We relied on the computed AOA surface to enforce predictor space similarity with the training data, thus ensuring that our model did not predict into unknown predictor space, even in coverages with limited FIA plots. Moreover, the AOA surface provided evidence of predictorspace uniformity across all 17 component LiDAR coverages, indicating that each of the component coverages was well represented in the model training dataset (Meyer and Pebesma, 2021). It stands to reason, however, that the NYC coverage contained the lowest proportion of AOA, since there were only two model plots available in this coverage (Table 1). Generally, pixels falling outside the AOA surface appeared to be the result of problems with LiDAR collection or data processing abnormalities, with some visible outliers that could not be attributed to any known ecological phenomena. We found the AOA mapping especially valuable in utilizing publicly available LiDAR coverages off the shelf with limited knowledge of, or responsibility for, their provenance.

4.3. Forest definition disparities

A forest can be defined in many ways depending on goals, perspectives, and operational concerns, and aligning the various definitions to make comparisons or derive relationships is neither a trivial nor a unique challenge (Chazdon et al., 2016; Riemann et al., 2010; Huang et al., 2019). FIA's strict forest definition that is in part based on field observations of land-use, which is traditionally difficult to classify with remotely sensed data (Fritz et al., 2017), was difficult to harmonize with LiDAR. Since FIA does not provide a forest/nonforest map, and FIA's definition of forest can exclude significant AGB stocks in areas containing tree cover (Johnson et al., 2015, 2014), we relied on LCMAP's classifications of vegetated cover types to mask our prediction surfaces. Our AGB maps thus reflect a more inclusive definition of forest than FIA, incorporating AGB stocks across a broader range of conditions and land-uses. Despite these definitional differences, there was overlap between FIA forest and our LCMAP-derived definition, as evidenced by the non-zero FIA AGB averages for model dataset plots grouped within each of the four vegetated LCMAP classes (Table 5). Further, we were able to separate true zeros from structural zeros in FIA nonforest conditions using a 1 m LiDAR max-height threshold (Table 2), giving our model information to make predictions in areas with little to no canopy cover.

Although most of our mapped AGB was contained within pixels classified as tree cover, a significant minority was contained within pixels classified as Cropland and Grass/Shrub (Table 5). As the total area of agricultural lands has been declining in NYS (USDA National Agricultural Statistics Service, 2019), we expect that many of these predictions represented early successional forests, or patches of 'young and stunted trees' in open fields (Yang et al., 2018), both transitional states which are challenges for LCMAP's classification algorithm (Mahoney et al., 2022; Brown et al., 2020). In summary, the landscape-level context in NYS, and LCMAP's algorithmic challenges provided support to generate predictions for all LCMAP vegetated classes.

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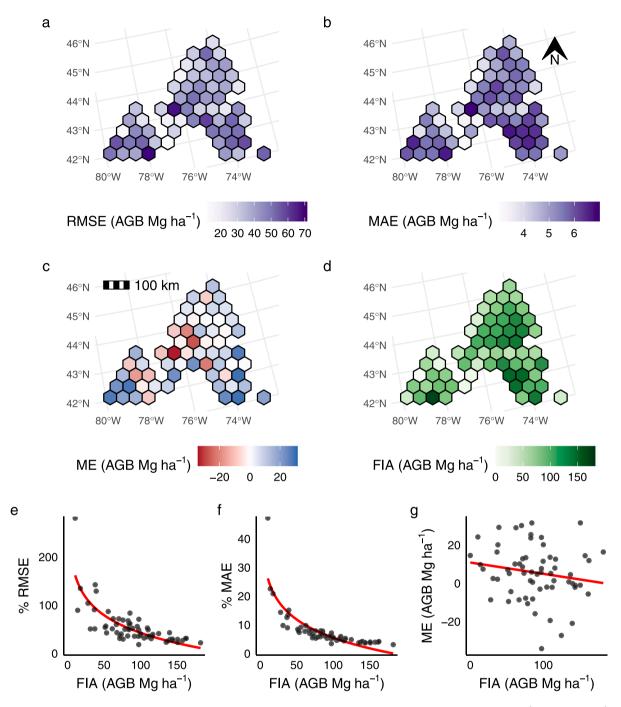


Fig. 7. Plot-to-pixel residuals summarized at units spaced 50 km apart. Hexagons with only one reference plot were removed. (a) RMSE Mg ha⁻¹; (b) MAE Mg ha⁻¹; (c) ME Mg ha⁻¹; (d) Mean FIA AGB value Mg ha⁻¹; (e) Hex-level % RMSE as a function of mean reference value; (f) Hex-level % MAE as a function of mean reference value; (g) Hex-level ME as a function of mean reference value; (f) where 0 FIA AGB makes % RMSE impossible to compute. Trend lines in (e) and (f) produced using logarithmic regression. Trend line in g produced using least-squares regression. RMSE, MAE, and ME as defined in Section 2.5.

4.4. Model performance and map agreement assessment

Our prediction accuracy against the 20% testing partition of our model dataset was favorably comparable to previous LiDAR-AGB mapping studies (Huang et al., 2019; Nilsson et al., 2017; Ayrey et al., 2021; Hauglin et al., 2021). Using a set of FIA-developed methods (Riemann et al., 2010; Menlove and Healey, 2020) we further demonstrated an overall strong agreement between our map-based estimates and FIA-derived estimates.

It is unsurprising that assessment metrics generally improved as the scales of aggregation increased, given that the plot-to-pixel scale can be considered the most rigorous, with the largest variance in both the pixel and plot AGB distributions, as well as the most potential for spatial misalignment between 30 m pixel predictions and FIA plot measurements to influence agreement metrics (McRoberts et al., 2018). When mapped residuals were summarized within units spaced 50 km apart, larger magnitudes of prediction error (RMSE, ME) emerged, but region or coverage specific patterns were not evident (Fig. 7). Rather, we observed ME to be mostly positive across the GPO-LiDAR area, and to be weakly related to the underlying distribution of FIA reference data. We can also infer that large RMSE values were a reflection of extreme individual outliers as they were often paired with reasonable MAE values.

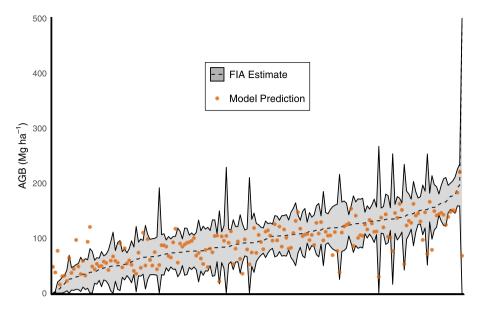


Fig. 8. Comparison of mapped predicted AGB to Menlove and Healey (2020) estimates (dashed line) and associated 95% confidence interval (gray shaded region) within 64,000-ha aggregation hexagons. FIA estimates of AGB are scaled by the proportion of forest cover indicated by LCMAP 2016 Tree cover, Wetland, Croplands, and Grass/Shrub classified pixels. Hexagons with mapped areas \leq 10% of their total area were excluded from this analysis. Observations are sorted by increasing FIA estimates along the *x*-axis.

Our choice to maintain FIA's probability sample in order to leverage unbiased estimators of map agreement metrics came with some drawbacks. Namely, we had to accept temporal lags of +/-2 years between LiDAR acquisitions and field inventories, and we had to include structural zeros in our assessment dataset where FIA did not record tree measurements in nonforest conditions. It is likely that the former inflated our estimates of error due to growth or disturbances occurring at plots between the time of inventory and the time of LiDAR acquisition, though we have no means to quantitatively confirm this possibility. We can say with more certainty that the structural zeros introduced by nonforest plots had a large impact on our map agreement assessment as evidenced by the stack of non-zero predictions along the x-axis in plot to pixel comparisons (Fig. 5) and the large positive ME for plots classified as Grass/Shrub and Cropland (Table 7). Given the large non-zero biomass predictions produced at these plots, it is not unreasonable to assume they contain trees unmeasured by FIA that would be captured by LiDAR height metrics. These structural zeros are likely the driving force behind the positive average prediction error (ME) found in our assessments, and may be the cause for the slight increase in Moran's I values computed with the assessment dataset relative to those computed with the model dataset (Supplementary Materials S7).

Nearly all of our assessments, including performance against the testing partition of the model data (Fig. 3), the Riemann analysis (Fig. 5), agreement with the Menlove and Healey FIA estimates (Fig. 8), and the weak relationship between ME and average FIA AGB estimate (Fig. 7g) reinforced our model's tendency towards lower accuracy at the extremes of the FIA reference dataset. These discrepancies can likely be attributed to the model structure and training approach, the aforementioned structural zeros in our map assessment dataset, as well as the saturation problem inherent in LiDAR-AGB modeling (St-Onge et al., 2008) where models fail to predict the largest AGB values in the data set. In general, with the extremes of the response distribution occurring less often, most models will be more accurate making predictions near the mean.

We also recognize the presence of uncertainty in the AGB reference data due to allometric, measurement, and locational errors, as well as our growth adjustment procedure used to boost the size of the model dataset, though quantifying the magnitude of this uncertainty was outside the scope of this paper. Duncanson et al. (2017) indicated that different choices in allometric models used to predict AGB from tree diameter and height measurements can result in large variation (up to 20%) of plot-level AGB estimates. An improvement would be to embed measurement errors, as well as allometric and growth adjustment uncertainty in a quantification of model precision or uncertainty at the pixel level (CEOS, 2021). Such information could help to develop variance estimates for any aggregation of pixel predictions that could be used in estimating AGB in areas that lack field inventories (Dettmann et al., 2022; CEOS, 2021).

4.5. Map applications

Our rigorously evaluated map products have a range of applications where knowledge of the spatial patterns of forest biomass (and by extension, forest carbon pools) is needed for monitoring, reporting, and verification efforts alongside policy or regulatory decision support. These uses include the identification of forested areas for future monitoring, protection, or management, and for providing AGB as a predictor in subsequent ecological models. Our AGB maps can also be leveraged as diverse training data for models driven by spaceborne remote sensing platforms with more contiguous spatial and temporal coverage than LiDAR patchworks, providing the basis for landscape-scale carbon accounting (Hudak et al., 2020; CEOS, 2021).

5. Conclusion

Accurate AGB predictions at fine resolutions can provide landowners and decision makers with valuable information on landscape patterns needed to implement forest-based climate solutions, including reforestation opportunities, avoided deforestation, and improved management for carbon storage and sequestration. We implemented a model-based approach leveraging extensive field inventory data (FIA) and publicly available LiDAR coverages to develop AGB maps across NYS, where forests are expected to contribute substantially as carbon sinks towards achieving a net-zero carbon economy by 2050. Although LiDAR point clouds provide detailed information on forest structure that can yield superior models of forest biomass, their limited coverage in both spatial and temporal domains produces patchworks of disparate datasets over broad scales. Our modeling approach, and the comprehensive set of assessments demonstrated here, addressed several of the common challenges inherent in using LiDAR patchworks for AGB mapping, including a lack of temporally matching reference data and data discrepancies among component LiDAR coverages. Our results show that our approach and the resulting map products provide accurate AGB information at scales relevant to forest and climate stewardship in NYS.

Declaration of competing interest

The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

Data availability

Data will be made available on request.

Acknowledgments

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Appendix A. Supplementary material

Supplementary material related to this article can be found online at https://doi.org/10.1016/j.jag.2022.103059.

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