

<u>Understanding Gradient</u> <u>Boosting as a gradient descent</u>

- For a given sample x_i , the predictions by a gradient-boosting regressor are given by $\Rightarrow \hat{y}_i = \sum_{m=1}^{n_{iter}} h_m(x_i) \text{ where } h_m \text{ is an instance of the base estimator}$
- It is often called a weak learner as it does not need to be extremely accurate
- It is mostly a decision tree and hence gradient boosting is generally performed using gradient boosting decision tree\
- The summation of all the base estimators gives the total gradient descent and each base estimator finds the gradient in its way in a function space unlike other algorithms where the most optimal parameter values are defined
- Gradient descent in parameter space: linear regression
 - The lease square loss is given by

$$egin{aligned} &\Rightarrow \mathcal{L} = \sum_i \left(\mathbf{y_i} - \hat{\mathbf{y_i}}
ight)^2 \ &\Rightarrow \mathcal{L} = \sum_i \left(\mathbf{y_i} - \mathbf{x_i^T} heta
ight)^2 ext{ where } \hat{\mathbf{y_i}} = \mathbf{x_i^T} heta \end{aligned}$$

The parameters are updated using the following

$$\Rightarrow$$
 $heta^{(\mathrm{m}\,+\,1)} = heta^{(\mathrm{m})} - lpha rac{\partial \mathcal{L}}{\partial heta^{(\mathrm{m})}}$

 $\circ~$ The optimal θ is the summation of all the terms where each term is the product of the learning rate and the negative gradient of the loss function

$$\Rightarrow \theta^{(m+2)} = \theta^{(m+1)} - \alpha \tfrac{\partial \mathcal{L}}{\partial \theta^{(m+1)}} = (\theta^{(m)} - \alpha \tfrac{\partial \mathcal{L}}{\partial \theta^{(m)}}) - \alpha \tfrac{\partial \mathcal{L}}{\partial \theta^{(m+1)}} \text{ using the power of induction}$$

- $\circ~$ In gradient boosting, the gradient is performed with respect to the predictions $\hat{y_i}$
- \circ Since gradient descent is used to minimise the loss function with respect to any of the independent variables, the predictions $\hat{y_i}$ which is also an independent variable, can be used to minimise the loss function

- · Minimising the loss with respect to the predictions
 - Predictions are randomly generated
 - The independent variables are updated using the following

$$\begin{split} &\Rightarrow \hat{\mathbf{y}_{\mathrm{i}}}^{(\mathrm{m}+1)} = \hat{\mathbf{y}_{\mathrm{i}}}^{(\mathrm{m})} - \alpha \frac{\partial \mathcal{L}}{\partial \hat{\mathbf{y}_{\mathrm{i}}}^{(\mathrm{m})}} \text{ where } \mathcal{L} = \sum_{i} \left(\mathbf{y}_{\mathrm{i}} - \hat{\mathbf{y}_{\mathrm{i}}}\right)^{2} \\ &\Rightarrow \frac{\partial \mathcal{L}}{\partial \hat{\mathbf{y}_{\mathrm{i}}}} = \frac{\partial}{\partial \hat{\mathbf{y}_{\mathrm{i}}}} \left(\mathbf{y}_{\mathrm{i}} - \hat{\mathbf{y}_{\mathrm{i}}}\right)^{2} \\ &\Rightarrow -2 \left(\mathbf{y}_{\mathrm{i}} - \hat{\mathbf{y}_{\mathrm{i}}}\right) \end{split}$$

- Gradient Boosting: use the base estimator to predict the gradients
 - Prediction takes place on unseen samples
 - Instead of updating the \hat{y} with the actual gradient, use a base estimator to predict the gradients
 - The set of base estimators can output a prediction for any sample
 - This can be represented by

$$\Rightarrow \hat{y}_i = \sum_{m \,=\, 1}^{n_{iter}} h_m(x_i)$$
 where h_m is an instance of the base estimator

- Instead of using gradient descent to estimate a parameter (a vector in a finite-dimensional space), use gradient descent to estimate a function (a vector in an infinite dimensional space)
- Wrapping up
 - Instead of taking the derivative of the loss with respect to the parameter of a parametrised model, take the derivative of the loss with respect to the predictions
 - There is no notion of a parametrised model anymore, but can still minimise our loss on the training samples
 - At each iteration of the gradient boosting procedure, train a base estimator to predict the gradient descent step
 - Saving these base estimators in memory is what enables to output predictions for any future sample