Classification as a task has a discrete output space. So we are trying to predict one of n classes. Therefore, we are either correct or wrong. In case of regression, the goal is to find a value in a more continuous space and hence even the most accurate model may be off by a few decimals.

2. SGD basically takes a subset of the training data to train the model in each iteration. In case of GD, each value of the error is calculated using the entire dataset. GD provides the more accurate estimate of the gradient. We can assume the gradient to be contributed from each training data and hence more the data, more accurate the estimate of the gradient. But this could lead to overfitting and hence noise in GD is helpful.

3. It is difficult to train neural networks with low initializations. This is because the activation value drop quickly to 0 when going into deeper layers. Therefore, during gradient descent, no learning happens as gradient is also 0.	s

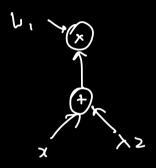
4. This is a bad idea. As all values are input and hence the same gradient de	equal, all the neurons in the hidden layer observe the same escent and in the end learn the same features.

$$\frac{\partial f}{\partial g} = \begin{cases} \alpha', & g < 0 \\ 1, & g > 0 \end{cases}$$

$$(I(g c_0). \propto + I(g >_0))$$

$$\frac{q\rho}{g\Gamma} = \frac{g\nu}{g\Gamma} = \frac{g\sigma}{g\Gamma} = (I(dco) \cdot \alpha + I(d>0)) (1-A)$$





W, update change

x does n'i

(x1>8) instead of

(x)

a = W, (x+>3)

OL 28 if 0 then same

- 3) a) hyperparameter -> Validation date.

 Not too large or small.
 - b) Not always, False Nevy high
 - c) Droport acts as ensembling. The model trains only
 on few inputs So do each overtite
 - a)a) log(0.3)
 - b) log (0.4)
 - C) (priect NOS more error.

Intention meximize the value of correct.

Not max amongst other classes.

d) How do you train?

- e) A -> Betch

 B-- Minibetch
- f) change in learning rate
- g) 720 ~0
- 3)a) Label
 - 6) he can one can with transfer territor