# Hydrus

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# Estimating Gaussians

In distributed systems

#### The Normal Distribution

$$\mathcal{N}(\mu, \sigma^2)$$

The Normal Distribution – Reparametrized

$$\mathcal{N}(\mu, \sigma^2) = \mathcal{N}\left(\mu, \frac{M_2}{n}\right)$$

The Normal Distribution – Reparametrized

$$\mathcal{N}(\mu, \sigma^2) = \mathcal{N}(n, \mu, M_2)$$

• Given an estimated Normal distribution,  $\mathcal{N}_t = \mathcal{N} (n_t, \mu_t, M_{2,t})$ 

• And a new data point,  $x_{t+1}$ 

• Update the distribution,  $\mathcal{N}_{t+1} = \mathcal{N} (n_{t+1}, \mu_{t+1}, M_{2,t+1})$ 

$$n_{t+1} = n_t + 1$$

$$\mu_{t+1} = \mu_t + \frac{x_{t+1} + \mu_t}{n_{t+1}}$$

$$M_{2,t+1} = M_{2,t} + (x_{t+1} - \mu_t)(x_{t+1} - \mu_{t+1})$$

```
def welford(dist, x):
 (count, mean, m2) = dist
 new_count = count + 1
 new_mean = mean + (x - mean) / count
 new_m2 = m2 + (x - mean) * (x - new_mean)
 return (new_count, new_mean, new_m2)
```

- Pros:
  - Easy to implement
  - Optimal time complexity: O(n)
- Cons:
  - Not parallel

• Given a dataset,  $X_{AB}$ 

• Partitioned into two,  $X_A$  and  $X_B$ 

ullet With distributions for both,  $\mathcal{N}_{\!A}$  and  $\mathcal{N}_{\!B}$ 

• Compute the combined distribution,  $\mathcal{N}_{AB}$ 

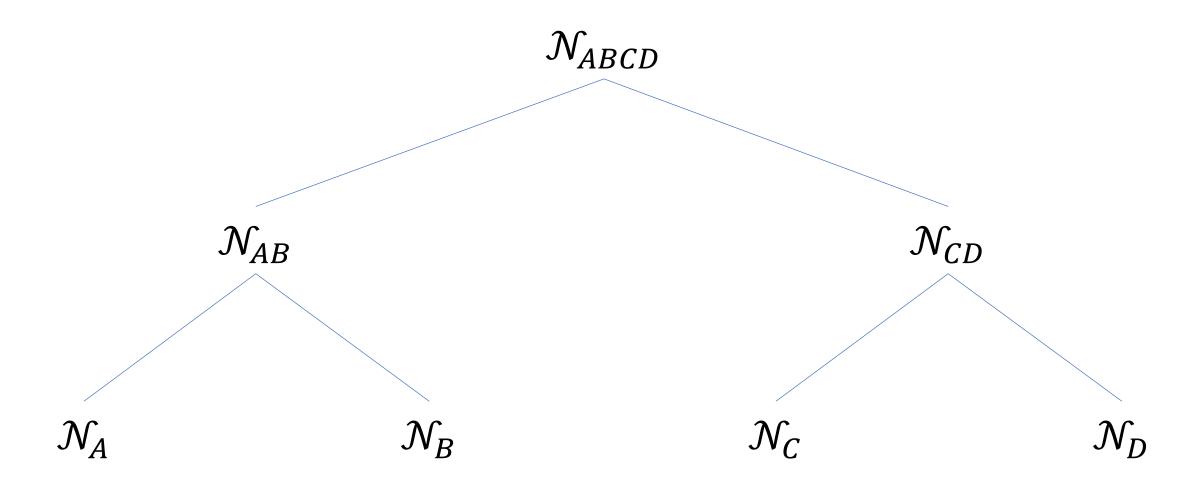
$$\delta = \mu_B - \mu_A$$

$$n_{AB} = n_A + n_B$$

$$\mu_{AB} = \mu_A + \delta \cdot \frac{n_B}{n_{AB}}$$

$$M_{2,AB} = M_{2,A} + M_{2,B} + \delta^2 \cdot \frac{n_A n_B}{n_{AB}}$$

```
def chan(dist a, dist b):
 (count a, mean a, m2 a) = dist a
 (count_b, mean_b, m2_b) = dist_b
delta = mean b - mean_a
 count = count_a + count_b
mean = mean a + delta * count b / count
m2 = m2 a + m2 b + delta**2
m2 *= count_a * count_b / count
 return (count, mean, m2)
```



- Pros
  - Easy to implement
  - Parallel
- Cons
  - Parallelism at the cost of CPU time
  - Possible stability issue with  $\mu_{AB}$ 
    - Alternate formula on Wikipedia

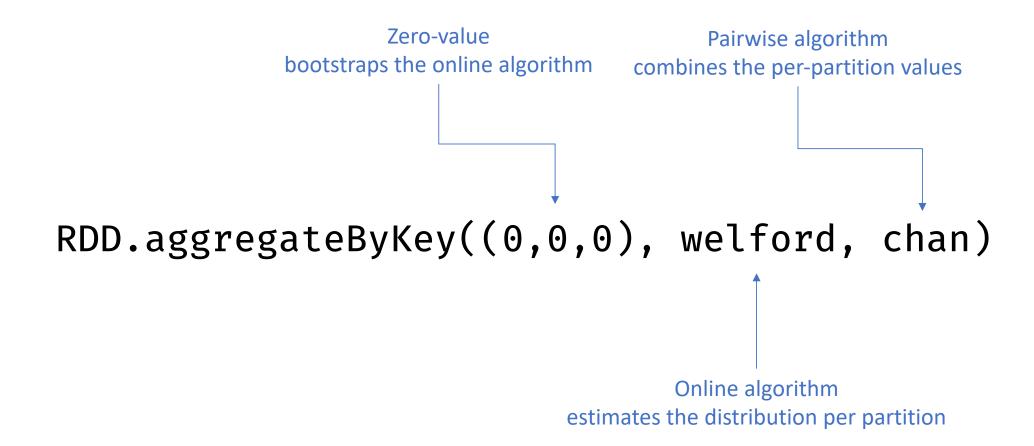
#### Spark

- Spark has methods like RDD.mean and RDD.variance
  - But not RDD.meanByKey or RDD.varianceByKey
  - Two passes over the data, one for each op
- We could use RDD.map and RDD.reduceByKey
  - Forces a shuffle
  - Easy to shoot your foot with a dumb initial key (n or  $M_2$ )
- Instead we have RDD.aggregateByKey

#### Spark

RDD.aggregateByKey((0,0,0), welford, chan)

#### Spark

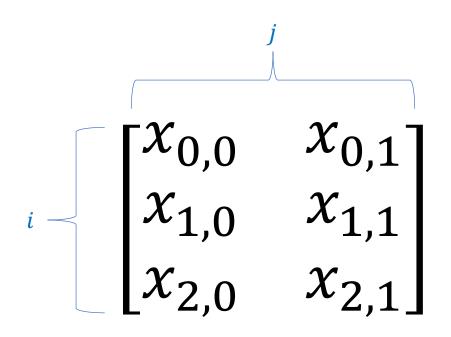




# Matrix Multiplication

with Spark RDDs

#### Representation



#### Representation

RDD[((
$$i, j$$
),  $x_{ij}$ )]

#### Representation

$$((i, j), x_{ij})$$

#### Matrix Multiplication

$$((i, j), x_{ij}) \times ((j, k), y_{jk})$$
  
=  $((i, k), \sum_{j} x_{ij} y_{jk})$ 

Step 0 – Start

$$((i, j), x_{ij}) \times ((j, k), y_{jk})$$

Step 1 – Rekey

$$(j, (i, x_{ij})) \times (j, (k, y_{jk}))$$

#### Step 2 – Join

$$(j, ((i, x_{ij}), (k, y_{jk})))$$

#### Step 3 – Multiply

$$(j, ((i, k), x_{ij}y_{jk}))$$

Step 4 – Drop j

$$((i, k), x_{ij}y_{jk})_{j}$$

Step 5 – Sum by key

$$((i, k), \sum_j x_{ij}y_{jk})$$

#### Step 6 – Profit!

- Ideal for sparse matrices
  - E.g. the document-term matrix
  - If the value is zero, don't include it in the RDD
  - The join step implicitly handles the zeros!
- *i*, *j*, and *k* don't need to be integers
  - Instead, use semantic values like document ID, word, and label

#### References

- Estimating Gaussians
  - https://en.wikipedia.org/wiki/Algorithms for calculating variance
  - B. P. Welford (1962),
     <u>"Note on a method for calculating corrected sums of squares and products"</u> (JSTOR),
     *Technometrics* 4(3):419–420
  - Chan, Tony F.; Golub, Gene H.; LeVeque, Randall J. (1979),
     <u>"Updating Formulae and a Pairwise Algorithm for Computing Sample Variances."</u> (PDF),
     Technical Report STAN-CS-79-773, Department of Computer Science, Stanford University.

