1) EM Algorithm

1) observed data vector:
$$X = (X_1, ..., X_n)$$
latent variable vector: $Z = (Z_1, ..., Z_n)$, each $Z_i = 0$ or 1
dishibution of Z : $X_i \mid Z_i = 0 \sim Pois(\mu_0)$, $Y_i \mid Z_i = 1 \sim Pois(\mu_i)$

$$P(Z_i = 0) = T$$
, $P(Z_i = 1) = 1 - TT$

2) Estep: solve for
$$Q(\theta | \theta^{(t)}) = estimating \theta$$
 bosed at $\theta^{(t)}$ from the t in iter, $Q(\theta | \theta^{(t)}) = E_{Z|X,\theta^{(t)}}(| \log L(\theta | X, Z)) = E_{Z|X,\theta^{(t)}}(| \sum_{i=1}^{\infty} \log L(\theta | X, Z))$ where $\theta = (\mu_i, \mu_i, \pi_i)$

$$= \sum_{i=1}^{\infty} \sum_{j=1}^{\infty} P(Z_{i} = j | X_{i} = x_{i}, \theta^{(t)}) \log L(\theta_{i} | X_{i}, j) \Rightarrow L(\theta_{i} | x_{i}, j) = \frac{\pi_{i} p_{\mu_{i}}(x_{i})}{\text{Top_{\mu_{0}}(X_{i})} - \pi_{i} p_{\mu_{i}}(x_{i})}$$

$$= \sum_{i=1}^{\infty} \sum_{j=1}^{\infty} T_{j,i}(| \log(\pi_{j} p_{\mu_{i}}(x_{i})) - \log(\pi_{0} p_{\mu_{0}}(x_{i}) + \pi_{i} p_{\mu_{i}}(x_{i}))), \text{ where } \pi_{0} = \pi_{0}$$

here
$$T_{j,i}^{(t)} := P(Z_{i=j}|X_{i=X_{i}}, M_{0}^{(t)}, M_{i}^{(t)}) = \frac{T_{i}^{(t)}P_{i}M_{i}^{(t)}(x_{i})}{T_{i}^{(t)}P_{i}M_{i}^{(t)}(x_{i})} = \frac{T_{i}^{(t)}P_{i}M_{i}^{(t)}(x_{i})}{T_{i}^{(t)}P_{i}M_{i}^{(t)}(x_{i})} = \frac{1}{N} \sum_{i=1}^{N} T_{i}(t_{i}) =$$

4) Use the result of k-means as an initial ostimator
$$\pi^{(0)} = \frac{\# \text{ of clustered 0}}{n}, \quad \Lambda_0^{(0)} = \frac{\text{Exic clustered 0}}{\# \text{ clustered 0}}, \quad \Lambda_1^{(0)} = \frac{\text{Exic clustered 1}}{\# \text{ clustered 1}}$$

Homework 3

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10/27/2021

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```
e_step <- function(x, z, pi, mu_0, mu_1) {</pre>
  # Input: x is the observation vector
           z is the hidden variable (0 or 1)
           pi, mu_0, mu_1 are the current parameters esitmates
  # Output: vector containing T_zi's
  \# PO(x) and P1(x) with current parameter estimates
  p0_x <- dpois(x, lambda = mu_0)</pre>
  p1_x <- dpois(x, lambda = mu_1)</pre>
  # compute the term in the denominator
  denom \leftarrow pi * p0_x + (1 - pi) * p1_x
  # compute the term in the numerator
  if (z == 0)
    num <- pi * p0_x
  else
    num <- (1 - pi) * p1_x
  return(num / denom)
m_step <- function(x, T_0, T_1) {</pre>
  # Input: x is the observation vector
         T_0 is nx1 vector of T_0i's
          T_1 is nx1 vector of T_1i's
  # Output: vector containing new estimates of pi, mu_0, mu_1
  mu_0 \leftarrow sum(T_0 * x) / sum(T_0)
  mu_1 \leftarrow sum(T_1 * x) / sum(T_1)
  pi \leftarrow mean(T_0)
  return(c(pi, mu_0, mu_1))
initial_vals <- function(x, n) {</pre>
  # Input: x is the observation vector
  # Output: vector containing initial guesses for pi, mu_0, mu_1
  kmeans <- kmeans(x, centers = 2)</pre>
```

```
# always label the first distribution as "1"
  if (kmeans$centers[1] > kmeans$centers[2]) {
    new_label = kmeans$cluster
    new label[kmeans$cluster == 1] = 2
    new label[kmeans$cluster == 2] = 1
    kmeans$cluster <- new_label</pre>
  pi <- sum(kmeans$cluster == 1) / n</pre>
  mu_0 <- mean(x[kmeans$cluster == 1])</pre>
  mu_1 <- mean(x[kmeans$cluster == 2])</pre>
  return(c(pi, mu_0, mu_1))
}
# return true if EM converged with threshold of 0.001
converged <- function(theta_prev, theta_curr) {</pre>
  dist <- abs(theta_prev - theta_curr) / theta_prev</pre>
  return(all(dist < 0.001))
}
```

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```
EM <- function(n, mu_0, mu_1) {</pre>
  # Input: n is the number of observations from each distribution
            so N in total is actually 2 * n
            mu_0, mu_1 are the true means used in simulations
  # Output: c(pi, mu_0, mu_1, accuracy), where accuracy is the
             assignment accuracy, and the first three elements are
             the estimates of the parameters from EM
  x_1 <- rpois(n, lambda = mu_0)</pre>
  x_2 \leftarrow rpois(n, lambda = mu_1)
  x \leftarrow c(x_1, x_2)
  init <- initial_vals(x, 2*n)</pre>
  theta_prev <- c(pi, mu_0, mu_1)</pre>
  T_0 \leftarrow e_{step}(x, 0, pi, mu_0, mu_1)
  T_1 \leftarrow e_{step}(x, 1, pi, mu_0, mu_1)
  theta_curr <- m_step(x, T_0, T_1)</pre>
  for (i in 1:5000) {
    if (converged(theta_prev, theta_curr)) break
    theta_prev <- theta_curr</pre>
    T_0 \leftarrow e_{step}(x, 0, theta_{curr}[1], theta_{curr}[2], theta_{curr}[3])
    T_1 <- e_step(x, 1, theta_curr[1], theta_curr[2],theta_curr[3])</pre>
    theta_curr <- m_step(x, T_0, T_1)
  result <- theta_curr
  T_0 \leftarrow e_{step}(x, 0, pi, mu_0, mu_1)
```

```
cluster <- rep(0, length(x))</pre>
  cluster [T_0 < .5] < 1  # assign i to cluster 0 if T_0 > .5
  # measure of accuracy
  cluster_true <- c(rep(0, n), rep(1, n)) # true clusters</pre>
  accuracy <- sum(cluster == cluster_true) / (2*n)</pre>
  result <- c(result, accuracy)</pre>
  return(result)
}
p6_1 \leftarrow EM(50, 5, 25)
p6_2 \leftarrow EM(50, 50, 10)
p6_3 <- EM(50, 100, 200)
p6_4 \leftarrow EM(50, 0.5, 5)
p6_5 \leftarrow EM(50, 40, 40)
acc \leftarrow c(p6_1[4], p6_2[4], p6_3[4], p6_4[4], p6_5[4])
barplot(acc, names.arg=c("(5, 25)", "(50, 10)", "(100, 200)", "(0.5, 5)", "(40, 40)"), xlab="(mu_0, mu_
               0.8
               9.0
               0.2
               0.0
                         (5, 25)
                                       (50, 10)
                                                    (100, 200)
                                                                    (0.5, 5)
                                                                                  (40, 40)
                                                  (mu_0, mu_1)
```

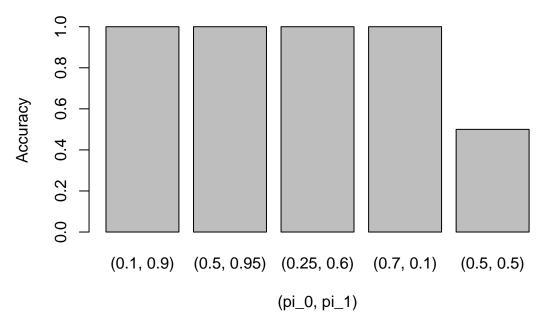
Bar chart shows that the accuracy of EM clustering with poisson distributions is very high. As expected, accuracy drops to 50% when the mixing distributions are identical (it's like choosing 50/50 every time).

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```
EM2 <- function(n, pi_0, pi_1) {
    # Input: n is the number of observations from each distribution
    # so N in total is actually 2 * n
    # Output: c(pi, mu_0, mu_1, accuracy), where accuracy is the
    # assignment accuracy, and the first three elements are
    # the estimates of the parameters from EM

x_1 <- rbinom(n, size = 100, prob = pi_0)
    x_2 <- rbinom(n, size = 100, prob = pi_1)
    x <- c(x_1, x_2)</pre>
```

```
mu_0 <- 100 * pi_0
  mu_1 <- 100 * pi_1
  init <- initial_vals(x, 2*n)</pre>
  theta_prev <- c(pi, mu_0, mu_1)</pre>
  T_0 \leftarrow e_{step}(x, 0, pi, mu_0, mu_1)
  T_1 <- e_step(x, 1, pi, mu_0, mu_1)</pre>
  theta_curr <- m_step(x, T_0, T_1)</pre>
  for (i in 1:5000) {
    if (converged(theta_prev, theta_curr)) break
    theta_prev <- theta_curr</pre>
    T_0 <- e_step(x, 0, theta_curr[1], theta_curr[2], theta_curr[3])</pre>
    T_1 <- e_step(x, 1, theta_curr[1], theta_curr[2],theta_curr[3])</pre>
    theta_curr <- m_step(x, T_0, T_1)
  result <- theta_curr
  T_0 \leftarrow e_{step}(x, 0, pi, mu_0, mu_1)
  cluster <- rep(0, length(x))</pre>
  cluster [T_0 < .5] < 1  # assign i to cluster 0 if T_0 > .5
  # measure of accuracy
  cluster_true <- c(rep(0, n), rep(1, n)) # true clusters</pre>
  accuracy <- sum(cluster == cluster_true) / (2 * n)</pre>
  result <- c(result, accuracy)</pre>
  return(result)
p7_1 \leftarrow EM2(50, 0.1, 0.9)
p7_2 \leftarrow EM2(50, 0.5, 0.95)
p7_3 \leftarrow EM2(50, 0.25, 0.6)
p7_4 \leftarrow EM2(50, 0.7, 0.1)
p7_5 \leftarrow EM2(50, 0.5, 0.5)
acc_7 \leftarrow c(p7_1[4], p7_2[4], p7_3[4], p7_4[4], p7_5[4])
barplot(acc_7, names.arg=c("(0.1, 0.9)", "(0.5, 0.95)", "(0.25, 0.6)", "(0.7, 0.1)", "(0.5, 0.5)"), xla
```



Similarly to previous results, this bar chart shows that the accuracy of poisson EM clustering with binomial distributions is very high. As expected, accuracy drops to 50% when the mixing distributions are identical (it's like choosing 50/50 every time).