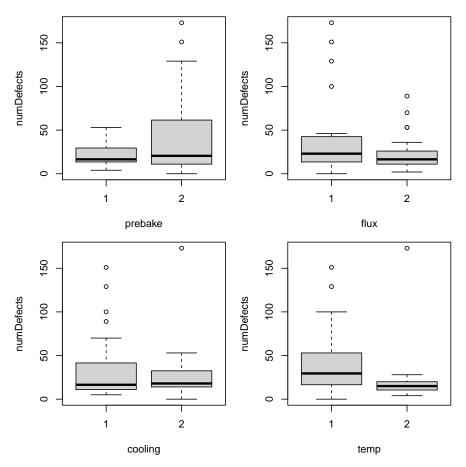
# MAST30027: Modern Applied Statistics

## Assignment 2 Solution 2021

1. Solution: There is no unique answer for this problem. This is a sample solution.

```
You can read the data using the following command.
```

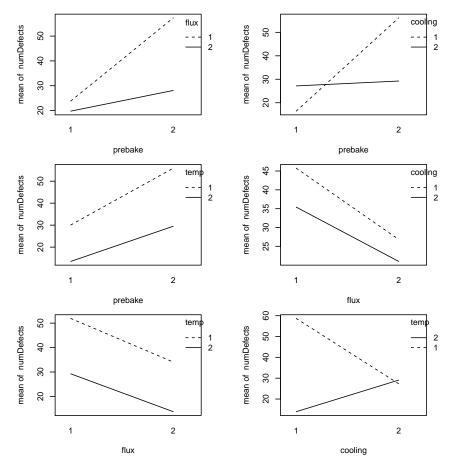
```
> data <- read.table(file ="assignment2_prob1_2021.txt", header=TRUE)</pre>
> dim(data)
[1] 48 5
> names(data)
[1] "numDefects" "prebake"
                                   "flux"
                                                  "cooling"
                                                                  "temp"
> data$prebake <- factor(data$prebake)</pre>
> data$flux <- factor(data$flux)</pre>
> data$cooling <- factor(data$cooling)</pre>
> data$temp <- factor(data$temp)</pre>
First, we visualize data by plotting the responses against the predictors.
(Predictors are factor, so we don't need to check linearity.)
> par(mfrow=c(2,2), mar=c(4,4,1,1))
> plot(numDefects ~ prebake, data)
> plot(numDefects ~ flux, data)
> plot(numDefects ~ cooling, data)
> plot(numDefects ~ temp, data)
```



The variable temp has a clear effect on the number of defects, but for the other variables things are less clear.

Are there any two way interactions between factors? We use interaction plots to check.

```
> par(mfrow=c(3,2))
> par(mar=c(4,4,1,4))
> with(data, interaction.plot(prebake,flux,numDefects))
> with(data, interaction.plot(prebake,cooling,numDefects))
> with(data, interaction.plot(prebake,temp,numDefects))
> with(data, interaction.plot(flux,cooling,numDefects))
> with(data, interaction.plot(flux,temp,numDefects))
> with(data, interaction.plot(cooling,temp,numDefects))
```



There are clear interactions between cooling and prebake, and between temp and cooling. Given that we have count data, a Poisson regression model with a log link is a good place to start. Since there are interactions, we use the main effects and all two-way interactions at once, then using stepwise selection with the AIC to select a parsimonious model.

```
> pmod <- glm(numDefects ~ (prebake + flux + cooling + temp)^2,
              family=poisson, data=data)
> pmod.f <- step(pmod, trace=0)</pre>
> summary(pmod.f)
Call:
glm(formula = numDefects ~ prebake + flux + cooling + temp +
    prebake:flux + prebake:cooling + prebake:temp + flux:temp +
    cooling:temp, family = poisson, data = data)
Deviance Residuals:
              1Q
                   Median
                                 ЗQ
                                         Max
                             1.3936
-7.7921
         -2.6541
                  -0.2946
                                    13.5042
Coefficients:
                  Estimate Std. Error z value Pr(>|z|)
(Intercept)
                               0.08598
                                       39.653
                                                < 2e-16 ***
                   3.40957
prebake2
                   1.33040
                               0.09625
                                        13.823
                                                < 2e-16 ***
flux2
                                                 0.3406
                  -0.09016
                               0.09461
                                        -0.953
cooling2
                   0.07550
                               0.09794
                                         0.771
                                                 0.4408
temp2
                  -1.86838
                               0.14883 -12.554 < 2e-16 ***
```

(Dispersion parameter for poisson family taken to be 1)

```
Null deviance: 1450.52 on 47 degrees of freedom Residual deviance: 626.99 on 38 degrees of freedom
```

AIC: 877.73

Number of Fisher Scoring iterations: 5

We see that the interaction term between flux and cooling was removed. We check the model adequacy.

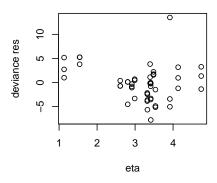
```
> pchisq(deviance(pmod.f), pmod.f$df.residual, lower.tail=FALSE)
```

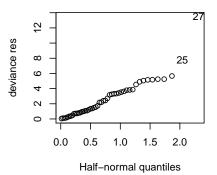
```
[1] 1.005731e-107
```

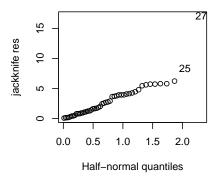
The p-value is very small, thus something is amiss with the model. The diagnostic plots might be able to tell us what is wrong.

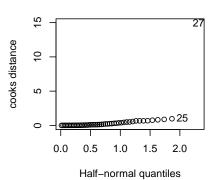
```
> library(faraway)
```

- > par(mfrow=c(2,2))
- > plot(predict(pmod.f), residuals(pmod.f), xlab="eta", ylab="deviance res")
- > halfnorm(residuals(pmod.f), ylab="deviance res")
- > halfnorm(rstudent(pmod.f), ylab="jackknife res")
- > halfnorm(cooks.distance(pmod.f), ylab="cooks distance")









Point 27 has very large Cook's distance, jackknife residual, and deviance residual, suggesting that this point might be an influential point. Thus, we will fit the model without the point 27

```
> pmod2 <- glm(numDefects ~ (prebake + flux + cooling + temp)^2,
+ family=poisson, subset=-27, data=data)
> pmod2.f <- step(pmod2, trace=0)
> summary(pmod2.f)
```

#### Call:

```
glm(formula = numDefects ~ prebake + flux + cooling + temp +
    prebake:cooling + flux:cooling + flux:temp + cooling:temp,
    family = poisson, data = data, subset = -27)
```

## Deviance Residuals:

${ t Min}$	1Q	Median	3Q	Max
-6.5090	-1.5709	-0.5039	1.5787	4.2854

#### Coefficients:

	${\tt Estimate}$	Std. Error	z value	Pr(> z )	
(Intercept)	3.52381	0.07795	45.205	< 2e-16	***
prebake2	1.23003	0.08099	15.187	< 2e-16	***
flux2	-0.57083	0.07473	-7.638	2.20e-14	***
cooling2	0.07465	0.10984	0.680	0.497	
temp2	-1.51196	0.10010	-15.104	< 2e-16	***
prebake2:cooling2	-1.77525	0.12359	-14.364	< 2e-16	***
flux2:cooling2	0.45881	0.11527	3.980	6.88e-05	***

---

Signif. codes: 0 '\*\*\* 0.001 '\*\* 0.01 '\* 0.05 '.' 0.1 ' ' 1

(Dispersion parameter for poisson family taken to be 1)

Null deviance: 1137.32 on 46 degrees of freedom Residual deviance: 289.02 on 38 degrees of freedom

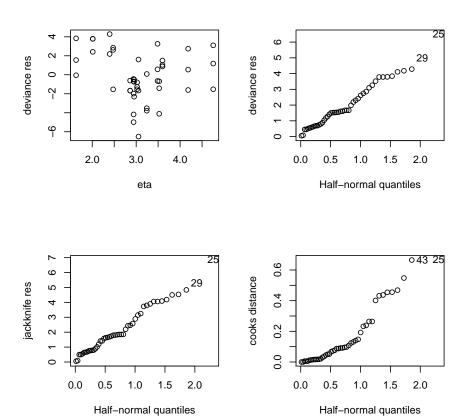
AIC: 530.77

Number of Fisher Scoring iterations: 5

> pchisq(deviance(pmod2.f), pmod2.f\$df.residual, lower.tail=FALSE)

## [1] 2.343993e-40

- > par(mfrow=c(2,2))
- > plot(predict(pmod2.f), residuals(pmod2.f), xlab="eta", ylab="deviance res")
- > halfnorm(residuals(pmod2.f), ylab="deviance res")
- > halfnorm(rstudent(pmod2.f), ylab="jackknife res")
- > halfnorm(cooks.distance(pmod2.f), ylab="cooks distance")



The diagnostic plots look fine now. However, we still observe that p-value of model adequacy test is low. Let's check overdispersion.

> (phi <- sum(residuals(pmod2.f, type="pearson")^2/pmod2.f\$df.residual))</pre>

#### [1] 7.355161

The large value of  $\hat{\phi}$  suggests that over dispersion occurs. Thus, we fit the data by quasi-Poisson model.

```
> pmod.q \leftarrow glm(numDefects \sim prebake + flux + cooling + temp + prebake:cooling + temp + temp + prebake:cooling + temp 
                               flux:cooling + flux:temp + cooling:temp,
                               family=quasipoisson, subset=-27, data=data)
+
> summary(pmod.q)
Call:
glm(formula = numDefects ~ prebake + flux + cooling + temp +
       prebake:cooling + flux:cooling + flux:temp + cooling:temp,
       family = quasipoisson, data = data, subset = -27)
Deviance Residuals:
       Min 1Q Median 3Q Max
-6.5090 -1.5709 -0.5039 1.5787
                                                                       4.2854
Coefficients:
                               Estimate Std. Error t value Pr(>|t|)
                                  (Intercept)
                                 0.21966 5.600 2.02e-06 ***
prebake2
flux2
cooling2
                          -1.51196 0.27148 -5.569 2.22e-06 ***
temp2
0.19085 0.34833 0.548 0.58696
flux2:temp2
Signif. codes: 0 '*** 0.001 '** 0.01 '* 0.05 '.' 0.1 ' 1
(Dispersion parameter for quasipoisson family taken to be 7.355163)
       Null deviance: 1137.32 on 46 degrees of freedom
Residual deviance: 289.02 on 38 degrees of freedom
AIC: NA
Number of Fisher Scoring iterations: 5
From the p-values in the summary table, flux2:cooling2 and flux2:temp2 are not signif-
icant. So we will consider a model without these two terms and compare the two models
using F-test.
> pmod.q.r <- glm(numDefects ~ prebake + flux + cooling + temp + prebake:cooling +
                                      cooling:temp, family=quasipoisson, subset=-27, data=data)
> anova(pmod.q.r, pmod.q, test="F")
Analysis of Deviance Table
Model 1: numDefects ~ prebake + flux + cooling + temp + prebake:cooling +
       cooling:temp
Model 2: numDefects ~ prebake + flux + cooling + temp + prebake:cooling +
        flux:cooling + flux:temp + cooling:temp
```

Resid. Df Resid. Dev Df Deviance F Pr(>F)

```
1 40 309.76
2 38 289.02 2 20.741 1.41 0.2566
```

The test shows that we should remove these two terms.

```
> summary(pmod.q.r)
```

#### Call:

```
glm(formula = numDefects ~ prebake + flux + cooling + temp +
    prebake:cooling + cooling:temp, family = quasipoisson, data = data,
    subset = -27)
```

#### Deviance Residuals:

```
Min 1Q Median 3Q Max
-6.893 -1.676 -0.090 1.432 4.104
```

#### Coefficients:

```
Estimate Std. Error t value Pr(>|t|)
                3.4384 0.2091 16.448 < 2e-16 ***
(Intercept)
prebake2
                 1.2300
                            0.2212 5.561 1.95e-06 ***
flux2
                 -0.3502
                            0.1494 -2.344 0.0241 *
                 0.2601
                            0.2675 0.972
                                            0.3367
cooling2
                            0.2351 -6.121 3.19e-07 ***
                 -1.4388
temp2
prebake2:cooling2 -1.7608
                            0.3375 -5.218 5.90e-06 ***
cooling2:temp2
                  0.8914
                            0.3472
                                    2.567 0.0141 *
```

Signif. codes: 0 '\*\*\*' 0.001 '\*\*' 0.05 '.' 0.1 ' ' 1

(Dispersion parameter for quasipoisson family taken to be 7.458044)

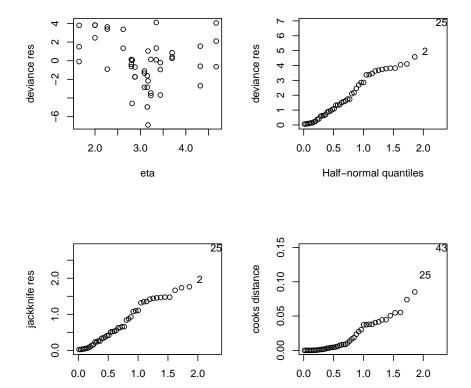
Null deviance: 1137.32 on 46 degrees of freedom Residual deviance: 309.76 on 40 degrees of freedom

AIC: NA

Number of Fisher Scoring iterations: 5

Now the summary looks fine. Note the main effect cooling is not significant. However, the interaction terms contains cooling. Hence, we will keep the main effect.

- > par(mfrow=c(2,2))
- > plot(predict(pmod.q.r), residuals(pmod.q.r), xlab="eta", ylab="deviance res")
- > halfnorm(residuals(pmod.q.r), ylab="deviance res")
- > halfnorm(rstudent(pmod.q.r), ylab="jackknife res")
- > halfnorm(cooks.distance(pmod.q.r), ylab="cooks distance")



The diagnostic plots of the resulting model also look fine.

Half-normal quantiles

flux is the only main effect which does not interact with any other factors. It has a negative effect on the number of defects since its coefficient is negative. prebake has a positive main effect on the number of defects. However, the coefficient of the interaction between prebake and cooling is negative and the effect is larger than the main effect of prebake. This means that the effect of prebake is negative when cooling = 2. (Other interpretations are acceptable as long as they are well explained.)

Half-normal quantiles

## 2. Solution:

(a) Let 
$$\theta = (\pi_1, \pi_2, \lambda_1, \lambda_2, \lambda_3)$$
 and  $n = 300$ . 
$$p(X_1, ...X_n, Z_1, ..., Z_n | \theta) = \prod_{i=1}^n p(X_i | Z_i, \theta) p(Z_i | \theta)$$
$$= \prod_{i=1}^n \prod_{k=1}^3 [p(X_i | Z_i = k, \theta) p(Z_i = k | \theta)]^{I_{(Z_i = k)}}.$$
$$\log[p(X_1, ...X_n, Z_1, ..., Z_n | \theta)] = \sum_{i=1}^n \sum_{k=1}^3 I_{(Z_i = k)} [\log p(X_i | Z_i = k, \theta) + \log p(Z_i = k | \theta)].$$

$$\begin{aligned} Q(\theta, \theta^0) &= E_{Z|X, \theta^0} [\log p(X_1, ... X_n, Z_1, ..., Z_n | \theta)] \\ &= \sum_{i=1}^n \sum_{k=1}^3 p(Z_i = k | X_i, \theta^0) [\log p(X_i | Z_i = k, \theta) + \log p(Z_i = k | \theta)] \\ &= \sum_{i=1}^n \sum_{k=1}^3 p(Z_i = k | X_i, \theta^0) [X_i \log \lambda_k - \lambda_k - \log X_i! + \log \pi_k], \end{aligned}$$

where  $\pi_3 = 1 - \pi_1 - \pi_2$ .

(b) Let  $\theta^0 = (\pi_1^0, \pi_2^0, \lambda_1^0, \lambda_2^0, \lambda_3^0)$ . E-step: For  $k \in \{1, 2\}$ ,

$$p(Z_i = k | X_i, \theta^0) = \frac{p(Z_i = k, X_i | \theta^0)}{p(X_i | \theta^0)}$$

$$= \frac{p(X_i | Z_i = k, \theta^0) p(Z_i = k | \theta^0)}{\sum_{s=1}^3 p(X_i | Z_i = s, \theta^0) p(Z_i = s | \theta^0)}$$

$$p(Z_i = 3 | X_i, \theta^0) = 1 - p(Z_i = 1 | X_i, \theta^0) - p(Z_i = 2 | X_i, \theta^0),$$

where  $p(X_i|Z_i=k,\theta^0) = \frac{(\lambda_k^0)^{X_i}e^{-\lambda_k^0}}{X_i!}$ ,  $p(Z_i=k|\theta^0) = \pi_k^0$  and  $p(Z_i=3|\theta^0) = 1 - \pi_1^0 - \pi_2^0$ 

(c) M-step:

$$\frac{\partial Q(\theta, \theta^0)}{\partial \pi_1} = \sum_{i=1}^n \left[ \frac{p(Z_i = 1 | X_i, \theta^0)}{\pi_1} - \frac{p(Z_i = 3 | X_i, \theta^0)}{1 - \pi_1 - \pi_2} \right] \\
= \frac{(1 - \pi_1 - \pi_2) \sum_{i=1}^n p(Z_i = 1 | X_i, \theta^0) - \pi_1 \sum_{i=1}^n p(Z_i = 3 | X_i, \theta^0)}{\pi_1 (1 - \pi_1 - \pi_2)} = 0 \quad (1)$$

$$\frac{\partial Q(\theta, \theta^0)}{\partial \pi_2} = \sum_{i=1}^n \left[ \frac{p(Z_i = 2|X_i, \theta^0)}{\pi_2} - \frac{p(Z_i = 3|X_i, \theta^0)}{1 - \pi_1 - \pi_2} \right] \\
= \frac{(1 - \pi_1 - \pi_2) \sum_{i=1}^n p(Z_i = 2|X_i, \theta^0) - \pi_2 \sum_{i=1}^n p(Z_i = 3|X_i, \theta^0)}{\pi_1 (1 - \pi_1 - \pi_2)} = 0. (2)$$

Let  $\pi_3 = 1 - \pi_1 - \pi_2$ . From (1) and (2), we obtain

$$\sum_{i=1}^{n} p(Z_i = 1 | X_i, \theta^0) \pi_3 = \sum_{i=1}^{n} p(Z_i = 3 | X_i, \theta^0) \pi_1,$$
(3)

$$\sum_{i=1}^{n} p(Z_i = 2|X_i, \theta^0) \pi_3 = \sum_{i=1}^{n} p(Z_i = 3|X_i, \theta^0) \pi_2.$$
(4)

Taking sum of (3) and (4), we have

$$\pi_3 \sum_{i=1}^n \left[ p(Z_i = 1 | X_i, \theta^0) + p(Z_i = 2 | X_i, \theta^0) \right] = (1 - \pi_3) \sum_{i=1}^n p(Z_i = 3 | X_i, \theta^0),$$
$$\hat{\pi}_3 = \frac{\sum_{i=1}^n p(Z_i = 3 | X_i, \theta^0)}{n}.$$

From (3), we have

$$\hat{\pi}_1 = \frac{1}{\sum_{i=1}^n p(Z_i = 3|X_i, \theta^0)} \hat{\pi}_3 \sum_{i=1}^n p(Z_i = 1|X_i, \theta^0) = \frac{\sum_{i=1}^n p(Z_i = 1|X_i, \theta^0)}{n}.$$

From (4), we have

$$\hat{\pi}_2 = \frac{1}{\sum_{i=1}^n p(Z_i = 3|X_i, \theta^0)} \hat{\pi}_3 \sum_{i}^n p(Z_i = 2|X_i, \theta^0) = \frac{\sum_{i=1}^n p(Z_i = 2|X_i, \theta^0)}{n}.$$

For k = 1, 2, 3,

$$\begin{split} \frac{\partial Q(\theta, \theta^0)}{\partial \lambda_k} &= \sum_{i=1}^n p(Z_i = k | X_i, \theta^0) \Big[ \frac{X_i}{\lambda_k} - 1 \Big] \\ &= \sum_{i=1}^n p(Z_i = k | X_i, \theta^0) \frac{X_i}{\lambda_k} - \sum_{i=1}^n p(Z_i = k | X_i, \theta^0) = 0. \\ \hat{\lambda}_k &= \frac{\sum_{i=1}^n p(Z_i = k | X_i, \theta^0) X_i}{\sum_{i=1}^n p(Z_i = k | X_i, \theta^0)}. \end{split}$$

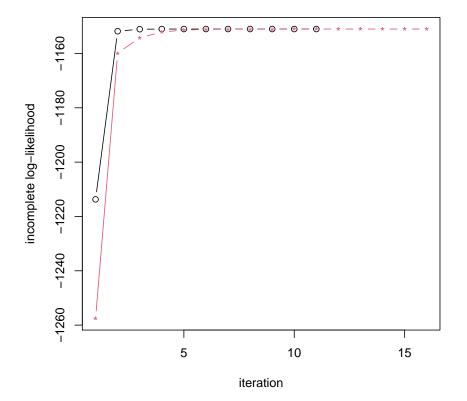
(d) Implement the EM algorithm.

```
> # w.init : initial value for pi
> # lambda.init : initial value for lambda
> # epsilon : stop if the change of the incomplete log-likelihood is less than epsilon
> # max.iter : maximum number of EM-iterations
> mixture.EM <- function(X, w.init, lambda.init, epsilon=1e-5, max.iter=100) {
    w.curr = w.init
   lambda.curr = lambda.init
    # store incomplete log-likehoods for each iteration
   log_liks = c()
    # compute incomplete log-likehoods using initial values of parameters.
    log_liks = c(log_liks, compute.log.lik(X, w.curr, lambda.curr)$ill)
    # set the change in incomplete log-likelihood with 1
   delta.ll = 1
    # number of iteration
   n.iter = 1
    # If the log-likelihood has changed by less than epsilon, EM will stop.
    while((delta.ll > epsilon) & (n.iter <= max.iter)){</pre>
+
      # run EM step
      EM.out = EM.iter(X, w.curr, lambda.curr)
+
      # replace the current value with the new parameter estimate
      w.curr = EM.out$w.new
      lambda.curr = EM.out$lambda.new
      # incomplete log-likehoods with new parameter estimate
      log_liks = c(log_liks, compute.log.lik(X, w.curr, lambda.curr)$ill)
      # compute the change in incomplete log-likelihood
      delta.ll = log_liks[length(log_liks)] - log_liks[length(log_liks)-1]
      # increase the number of iteration
```

```
n.iter = n.iter + 1
+
+
    }
+
    return(list(w.curr=w.curr, lambda.curr=lambda.curr, log_liks=log_liks))
+ }
> EM.iter <- function(X, w.curr, lambda.curr) {
    # E-step: compute E_{Z/X}, theta_0}[I(Z_i = k)]
+
+
    # for each sample X_i, compute P(X_i, Z_i=k)
    prob.x.z = compute.prob.x.z(X, w.curr, lambda.curr)$prob.x.z
    # compute P(Z_i=k \mid X_i)
    P_ik = prob.x.z / rowSums(prob.x.z)
    # M-step
    w.new = colSums(P_ik)/sum(P_ik) # sum(P_ik) is equivalent to sample size
    lambda.new = colSums(P_ik*X)/colSums(P_ik)
+
    return(list(w.new=w.new, lambda.new=lambda.new))
+ }
> # for each sample $X_i$, compute $P(X_i, Z_i=k)$
> compute.prob.x.z <- function(X, w.curr, lambda.curr) {</pre>
    # for each sample X_i, compute P(X_i, Z_i=k).
   # store these values in the columns of L:
   L = matrix(NA, nrow=length(X), ncol= length(w.curr))
   for(k in seq_len(ncol(L))) {
      L[, k] = dpois(X, lambda.curr[k])*w.curr[k]
    return(list(prob.x.z=L))
+
+ }
> # Compute incomplete log-likehoods
> compute.log.lik <- function(X, w.curr, lambda.curr) {</pre>
    # for each sample X_i, compute P(X_i, Z_i=k)
    prob.x.z = compute.prob.x.z(X, w.curr, lambda.curr)$prob.x.z
    # incomplete log-likehoods
    ill = sum(log(rowSums(prob.x.z)))
   return(list(ill=ill))
Run the EM algorithm two times with the two initial values provided in the problem.
> # read the data
> X = scan(file="assignment2_prob2_2021.txt", what=double())
> EM1 < -mixture.EM(X, w.init=c(0.3,0.3,0.4), lambda.init=c(3, 20, 35),
                    epsilon=1e-5, max.iter=100)
> EM2 <- mixture.EM(X, w.init=c(0.1,0.2,0.7), lambda.init=c(5, 25, 40),
                    epsilon=1e-5, max.iter=100)
> check = rbind(c(EM1$w.curr, EM1$lambda.curr),
```

Estimates from the two EM runs are very similar and they have very similar incomplete log-likelihoods. So I will report one from the second run as it has slightly higher incomplete log-likelihoods. MLEs for the parameters are  $\hat{\pi}_1 = 0.2491307, \hat{\pi}_2 = 0.2497786, \hat{\lambda}_1 = 5.167466, \hat{\lambda}_2 = 18.09381, \hat{\lambda}_3 = 36.93911.$ 

Check that the incomplete log-likelihoods increases at each step by plotting them



### 3. Solution:

(a) Let 
$$\theta = (\pi_1, \pi_2, \lambda_1, \lambda_2, \lambda_3), n = 300 \text{ and } n' = 100.$$

$$\begin{aligned} & p(X_1, ...X_n, Z_1, ..., Z_n, X_{n+1}, ..., X_{n+n'}|\theta) \\ &= \prod_{i=1}^n p(X_i|Z_i, \theta) p(Z_i|\theta) \prod_{i=n+1}^{n+n'} p(X_i|\theta) \\ &= \prod_{i=1}^n \prod_{k=1}^3 [p(X_i|Z_i = k, \theta) p(Z_i = k|\theta)]^{I(Z_i = k)} \prod_{i=n+1}^{n+n'} p(X_i|\theta). \end{aligned}$$

$$\log[p(X_1, ... X_n, Z_1, ..., Z_n, X_{n+1}, ..., X_{n+n'}|\theta)]$$

$$= \sum_{i=1}^n \sum_{k=1}^3 I_{(Z_i=k)}[\log p(X_i|Z_i=k, \theta) + \log p(Z_i=k|\theta)] + \sum_{i=n+1}^{n+n'} \log p(X_i|\theta)$$

$$\begin{split} Q(\theta,\theta^0) &= E_{Z|X,\theta^0}[\log p(X_1,...X_n,Z_1,...,Z_n,X_{n+1},...,X_{n+n'}|\theta)] \\ &= \sum_{i=1}^n \sum_{k=1}^3 p(Z_i = k|X_i,\theta^0)[\log p(X_i|Z_i = k,\theta) + \log p(Z_i = k|\theta)] + \sum_{i=n+1}^{n+n'} \log p(X_i|\theta) \\ &= \sum_{i=1}^n \sum_{k=1}^3 p(Z_i = k|X_i,\theta^0)[X_i\log \lambda_k - \lambda_k - \log X_i! + \log \pi_k] + \sum_{i=n+1}^{n+n'} [X_i\log \lambda_2 - \lambda_2 - \log X_i!], \end{split}$$

where  $\pi_3 = 1 - \pi_1 - \pi_2$ .

(b) E-step: same as the E-step in the solution for the problem 2 (b).

M-step: same as the M-step in the solution for the problem 2 (c) except for the following:

$$\frac{\partial Q(\theta, \theta^0)}{\partial \lambda_2} = \sum_{i=1}^n p(Z_i = 2|X_i, \theta^0) \left[ \frac{X_i}{\lambda_2} - 1 \right] + \sum_{i=n+1}^{n+n'} \left[ \frac{X_i}{\lambda_2} - 1 \right]$$
$$= \sum_{i=1}^n p(Z_i = 2|X_i, \theta^0) \frac{X_i}{\lambda_2} - \sum_{i=1}^n p(Z_i = 2|X_i, \theta^0) + \sum_{i=n+1}^{n+n'} \frac{X_i}{\lambda_2} - n' = 0.$$

Hence,

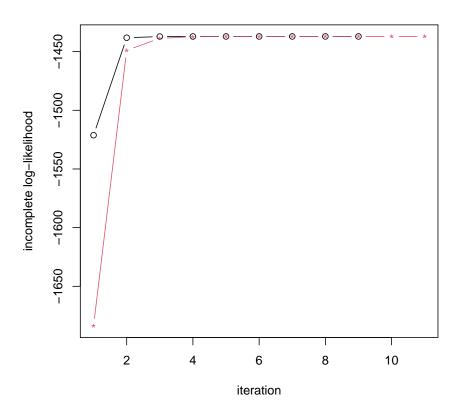
$$\hat{\lambda}_2 = \frac{\sum_{i=1}^n p(Z_i = 2|X_i, \theta^0) X_i + \sum_{i=n+1}^{n+n'} X_i}{\sum_{i=1}^n p(Z_i = 2|X_i, \theta^0) + n'}.$$

(c) Implement the EM algorithm.

> # X : X\_1, ..., X\_300 which follow a mixture of Poisson distribution
> # X0 : X\_301, ..., X\_400 which follow a Poisson distribution
> mixture.EM <- function(X, X0, w.init, lambda.init, epsilon=1e-5, max.iter=100) {
+
+ w.curr = w.init
+ lambda.curr = lambda.init
+
+ # store incomplete log-likehoods for each iteration
+ log\_liks = c()</pre>

```
# compute incomplete log-likehoods using initial values of parameters.
    log_liks = c(log_liks, compute.log.lik(X, X0, w.curr, lambda.curr)$ill)
    # set the change in incomplete log-likelihood with 1
    delta.ll = 1
    # number of iteration
    n.it.er = 1
    # If the log-likelihood has changed by less than epsilon, EM will stop.
    while((delta.ll > epsilon) & (n.iter <= max.iter)){</pre>
      # run EM step
      EM.out = EM.iter(X, X0, w.curr, lambda.curr)
      # replace the current value with the new parameter estimate
      w.curr = EM.out$w.new
+
      lambda.curr = EM.out$lambda.new
      # incomplete log-likehoods with new parameter estimate
      log_liks = c(log_liks, compute.log.lik(X, X0, w.curr, lambda.curr)$ill)
+
      # compute the change in incomplete log-likelihood
      delta.ll = log_liks[length(log_liks)] - log_liks[length(log_liks)-1]
      # increase the number of iteration
     n.iter = n.iter + 1
    return(list(w.curr=w.curr, lambda.curr=lambda.curr, log_liks=log_liks))
+
+ }
> EM.iter <- function(X, XO, w.curr, lambda.curr) {
+
    # E-step: compute E_{Z|X, \theta}[I(Z_i = k)]
    # for each sample X_i, compute P(X_i, Z_i=k)
    prob.x.z = compute.prob.x.z(X, w.curr, lambda.curr)$prob.x.z
    # compute P(Z_i=k \mid X_i)
   P_ik = prob.x.z / rowSums(prob.x.z)
    # M-step
   w.new = colSums(P_ik)/sum(P_ik) # sum(P_ik) is equivalent to sample size
    ### Change!!!
    lambda.new = rep(NA, length(w.new))
    lambda.new[c(1,3)] = (colSums(P_ik*X)/(colSums(P_ik)))[c(1,3)]
    lambda.new[2] = (colSums(P_ik*X)[2] + sum(XO))/(colSums(P_ik)[2] + length(XO))
+
    return(list(w.new=w.new, lambda.new=lambda.new))
+ }
> # for each sample X_i, compute P(X_i, Z_i=k)
> compute.prob.x.z <- function(X, w.curr, lambda.curr) {</pre>
    # for each sample X_i, compute P(X_i, Z_i=k).
```

```
# Store these values in the columns of L:
    L = matrix(NA, nrow=length(X), ncol= length(w.curr))
    for(k in seq_len(ncol(L))) {
      L[, k] = dpois(X, lambda.curr[k])*w.curr[k]
+
    return(list(prob.x.z=L))
+ }
> # Compute incomplete log-likehoods
> compute.log.lik <- function(X, X0, w.curr, lambda.curr) {</pre>
    # for each sample $X_i$, compute $P(X_i, Z_i=k)$
    prob.x.z = compute.prob.x.z(X, w.curr, lambda.curr)$prob.x.z
    # incomplete log-likehoods
    ill = sum(log(rowSums(prob.x.z))) + sum(log(dpois(X0, lambda.curr[2])))
+
    return(list(ill=ill))
+ }
Run the EM algorithm two times with the two initial values provided in the problem.
> X0 = scan(file="assignment2_prob3_2021.txt", what=double())
> EM1 < -mixture.EM(X, X0, w.init=c(0.3,0.3,0.4), lambda.init=c(3, 20, 35),
                     epsilon=1e-5, max.iter=100)
> EM2 <- mixture.EM(X, X0, w.init=c(0.1,0.2,0.7), lambda.init=c(5, 25, 40),
                     epsilon=1e-5, max.iter=100)
> check = rbind(c(EM1$w.curr, EM1$lambda.curr),
                 c(EM2$w.curr, EM2$lambda.curr))
> colnames(check) = c('pi_1', 'pi_2', 'pi_3', 'lambda_1', 'lambda_2', 'lambda_3')
> rownames(check) = c('EM1', 'EM2')
> check
                              pi_3 lambda_1 lambda_2 lambda_3
                    pi_2
         pi_1
EM1 0.2461850 0.2447837 0.5090313 5.118187 17.36453 36.76509
EM2 0.2461848 0.2447753 0.5090399 5.118186 17.36442 36.76488
> print(EM1$log_liks[length(EM1$log_liks)], digits=16)
[1] -1437.052845923898
> print(EM2$log_liks[length(EM2$log_liks)], digits=16)
[1] -1437.052845374368
Estimates from the two EM runs are very similar and they have very similar in-
complete log-likelihoods. So I will report one from the second run as it has slightly
higher incomplete log-likelihoods. MLEs for the parameters are \hat{\pi}_1 = 0.2461848, \hat{\pi}_2 =
0.2447753, \hat{\lambda}_1 = 5.118186, \hat{\lambda}_2 = 17.36442, \hat{\lambda}_3 = 36.76488.
Check that the incomplete log-likelihoods increases at each step by plotting them
> plot(EM1$log_liks, type='b',
       xlim=c(1,max(length(EM1$log_liks), length(EM2$log_liks))),
       ylim=range(c(EM1$log_liks, EM2$log_liks)),
       ylab='incomplete log-likelihood', xlab='iteration')
> points(EM2$log_liks, type='b', pch='*', col=2)
```



4. **NA.**