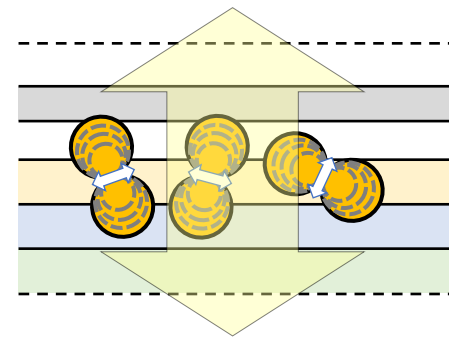


Internal Light Source- Power Density

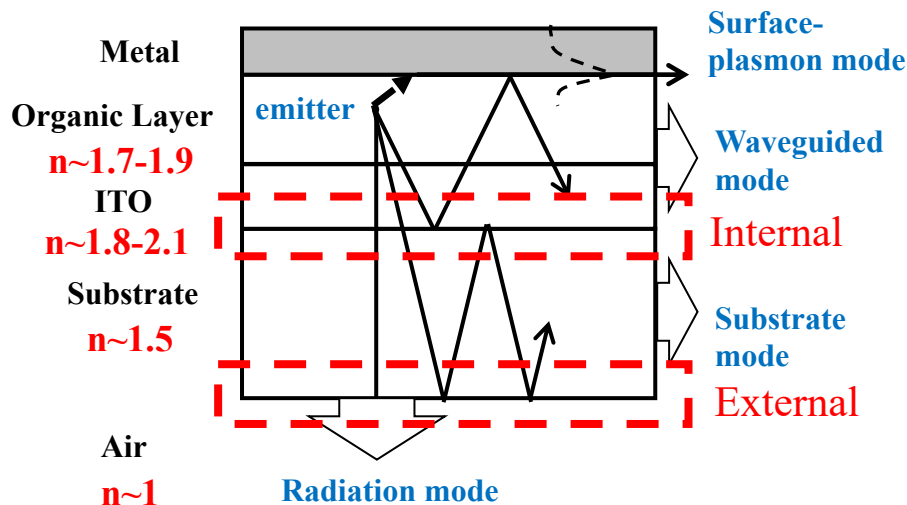
-PowerDensityCmd.pyc

Author: Wei-Kai Lee

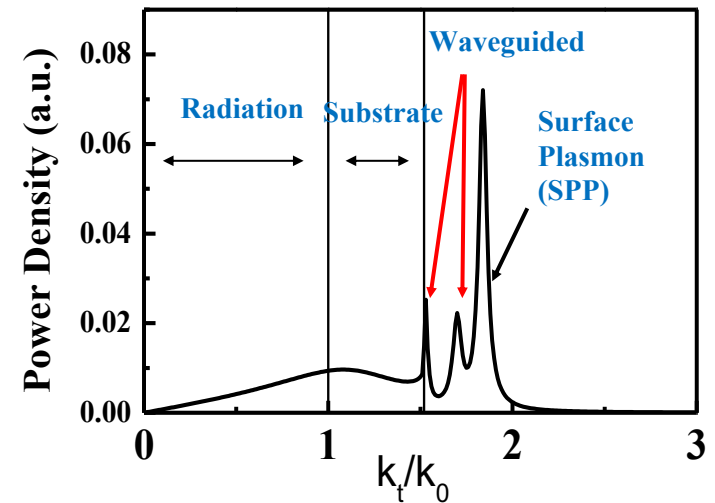


Objective

- To calculate the power density of the device.



Power Density/Power Dissipation



How to execute the calculator

python: windows
python3: mac, linux

```
h>python PowerDensityCmd.pyc
```

Execution file

```
>>> Please insert username : user-1
```

Type user name

```
Power Density Setting is not built in ../../SETTING/user-1  
Now saving the default power density setting.
```

LabSimulator_aniso > LegendDesign > SETTING > user-1	
名稱	修改日期
DFCmd.setting	2020/4/13 下
log	2020/4/14 上
materialMgr.mMgr	2020/4/13 下
PDCmd.setting	2020/4/14 上
rtauCmd.setting	2020/4/12 下
TRACmd.setting	2020/4/13 上

Help

```
>>> ?

User Control Command
=====
1. Setting Command:
changeUser      exit

Material Manager Command
=====
1. Setting Command:
printMgr        saveMgr

Structure/Structure List Command
=====
1. Structure List Command:
ReadStructListPath  ReadStructListName
SaveStructListPath  SaveStructListName
readStructList      saveStructList

2. Structure Command:
ReadStructPath      ReadStructName
SaveFilePath        SaveFileName
readStruct          deleteStruct

3. Print Information Command:
printStructInfo     printStructSettingInfo
printListInfo

4. Result Command:
ResultFilePath      ResultFileName
deleteResult
save_run_time_result_Bool
resetSN
```

Help

```
Power Density Command
**** Notice the mode interference power !!!
=====
1. Settting Command:
SettingFilePath      SettingFileName
setDefaultSetting    printPDInfo
loadPDSETTING        savePDSETTING

2. Power Density in kxky domain:
kxko                 kyko                 ----Parameter
runPD_kxky           plotPDvsKxky
plotPDkxkyContour    writekxkyMatrix

3. Power Density in k polar domain:
ktko                 Thetakt              ----Parameter
runPD_PolarInK       plotPDvsKpolar
plotPDPolarContour   writekPolarMatrix

4. Power Density z in kxky domain:
kxko                 kyko                 z                 ----Parameter
runPDz_kxky          plotPDvsZ_kxky

5. Power Density z in k polar domain:
ktko                 Thetakt              z                 ----Parameter
runPDz_PolarInK      plotPDvsZ_kPolar

6. Incoherence Layer Power Density in kxky domain:
kxko                 kyko                 ----Parameter
runInCoPD_kxky       plotInCoPDvsKxky
plotInCoPDkxkyContour writeInCokxkyMatrix

7. Incoherence Layer Power Density in k polar domain:
ktko                 Thetakt              ----Parameter
runInCoPD_PolarInK   plotInCoPDvsKpolar
plotInCoPDPolarContour writeInCokPolarMatrix
```

Help

```
8. Write Bool:
--- Mode (related to Power Density/ Power Density z) ---
write_TotalPowerBool
write_ForwardPowerBool      write_BackwardPowerBool
write_M1M3PowerBool         write_M2M4PowerBool
write_M1PowerBool           write_M2PowerBool
write_M3PowerBool           write_M4PowerBool
--- Region (Only related to Power Density) ---
write_EMLBool               ----at the emission position
write_EML_zPlusBool         ----at the position just above the emission position
write_EML_zMinusBool        ----at the position just below the emission position
write_SEMI_zPlusBool        ----at the position just above the 1st interface of the device
write_SEMI_zMinusBool       ----at the position just below the last interface of the device

9. Write Matrix Bool:
--- Mode (related to Power Density/ Power Density z/ Incoherence Power Density) ---
writeMatrix_TotalPowerBool
writeMatrix_ForwardPowerBool writeMatrix_BackwardPowerBool
writeMatrix_M1M3PowerBool    writeMatrix_M2M4PowerBool
writeMatrix_M1PowerBool      writeMatrix_M2PowerBool
writeMatrix_M3PowerBool      writeMatrix_M4PowerBool
--- Region (Only related to Power Density) ---
writeMatrix_EMLBool          ----at the emission position
writeMatrix_EML_zPlusBool    ----at the position just above the emission position
writeMatrix_EML_zMinusBool   ----at the position just below the emission position
writeMatrix_SEMI_zPlusBool   ----at the position just above the 1st interface of the device
writeMatrix_SEMI_zMinusBool  ----at the position just below the last interface of the device
```



Help

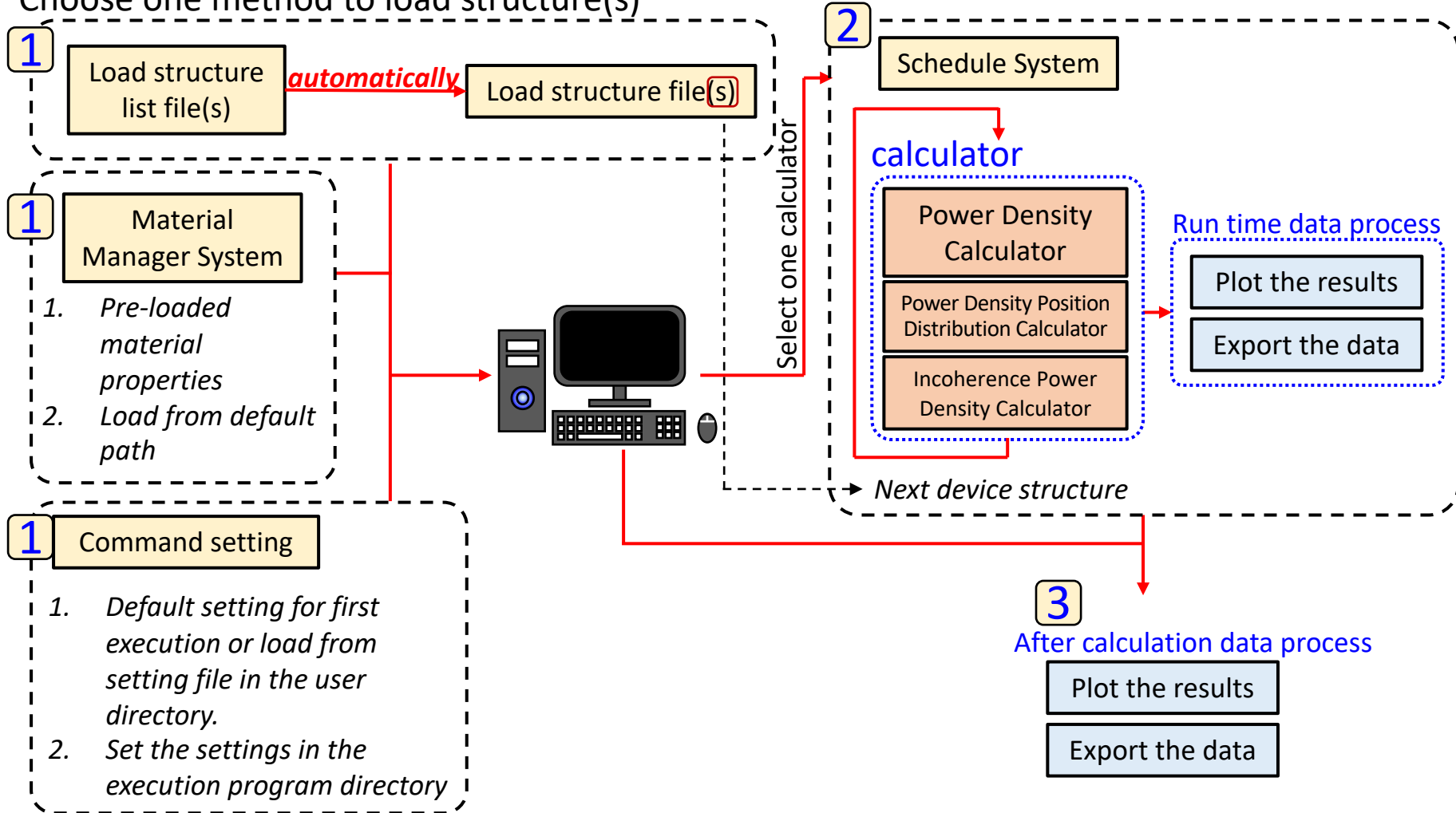
```
10. Plot Bool:
changeFigshowBool
--- Mode (related to Power Density/ Power Density z/ Incoherence Power Density) ---
plot_TotalPowerBool
plot_ForwardPowerBool      plot_BackwardPowerBool
plot_M1M3PowerBool         plot_M2M4PowerBool
plot_M1PowerBool           plot_M2PowerBool
plot_M3PowerBool           plot_M4PowerBool
--- Region (Only related to Power Density) ---
plot_EMLBool               ----at the emission position
plot_EML_zPlusBool         ----at the position just above the emission position
plot_EML_zMinusBool        ----at the position just below the emission position
plot_SEMI_zPlusBool        ----at the position just above the 1st interface of the device
plot_SEMI_zMinusBool       ----at the position just below the last interface of the device

11. Run Time Bool:
runtime_plot
--- (Only related to Power Density/ Incoherence Power Density)) ---
runtime_write_matrix runtime_contourplot
```



Calculating Workflow

Choose one method to load structure(s)



Default Setting

```
>>> printPDInfo

Setting file path : ../.. /SETTING/user-1
Setting file name : PDCmd.setting
kt/ko : 0.00000:0.10000:1.00000
Theta in kt/ko domain : 0.00000:10.00000:360.00000
kx/ko : -1.00000:0.10000:1.00000
ky/ko : -1.00000:0.10000:1.00000
z(nm) : -50.00000:1.00000:50.00000
Save Run Time Result : True
Run Time Write Matrix : False
Run Time Plot : False
Run Time Contour Plot : False

Plot Region Bool:
EML : True (EMLz+ add EMLz- )
EML(EMLz+) : False (at the position just above emitting position)
EML(EMLz-) : False (at the position just below emitting position)
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)
-Semi-infinite region: True (at the interface of the last semi-infinite region)
Plot Mode Bool: (should be cosidered with plot region bool)
Total (M1+M2+M3+M4) : True
+z (M1+M2) : False
-z (M3+M4) : False
M1 : False
M2 : False
M3 : False
M4 : False

Write Region Bool:
EML : True (EMLz+ add EMLz- )
EML(EMLz+) : True (at the position just above emitting position)
EML(EMLz-) : True (at the position just below emitting position)
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)
-Semi-infinite region: True (at the interface of the last semi-infinite region)
Write Mode Bool: (should be cosidered with plot region bool)
Total (M1+M2+M3+M4) : True
+z (M1+M2) : True
-z (M3+M4) : True
M1 : True
M2 : True
M3 : True
M4 : True

WriteMatrix Region Bool:
EML : True (EMLz+ add EMLz- )
EML(EMLz+) : False (at the position just above emitting position)
EML(EMLz-) : False (at the position just below emitting position)
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)
-Semi-infinite region: True (at the interface of the last semi-infinite region)
WriteMatrix Mode Bool: (should be cosidered with plot region bool)
Total (M1+M2+M3+M4) : True
+z (M1+M2) : False
-z (M3+M4) : False
M1 : False
M2 : False
```

Setting

```
Setting file name: PDCmdSetting  
kt/ko : 0.00000:0.10000:1.00000  
Theta in kt/ko domain : 0.00000:10.00000:360.00000
```

kt/ko and θ_{kt} setting method

```
<PDCmd> help ktco  
Set the kt/ko. (Empty for delete setting)  
[Usage] ktco [ktco] - single value, start:spacing:end, (v1,v2,v3,v4)  
  
<PDCmd> help Thetakt  
Set the polar angle in kx ky domain.  
[Usage] Thetakt [theta] - single value, start:spacing:end, (v1,v2,v3,v4)
```

Setting

```
kx/ko : -1.00000:0.10000:1.00000  
ky/ko : -1.00000:0.10000:1.00000
```

kx/ko and ky/ko setting method

```
<PDCmd> help kxko  
Set the kx/ko. (Empty for delete setting)  
[Usage] kxko [kxko] - single value, start:spacing:end, (v1,v2,v3,v4)  
  
<PDCmd> help kyko  
Set the ky/ko. (Empty for delete setting)  
[Usage] kyko [kyko] - single value, start:spacing:end, (v1,v2,v3,v4)
```

```
z(nm) : -50.00000:1.00000:50.00000
```

z setting method

```
<PDCmd> help z  
Set the simulation position z. (Empty for delete setting)  
[Usage] z [z] - single value, start:spacing:end, (v1,v2,v3,v4)
```

Run-Time Setting

```
Save Run Time Result : True
```

Whether to save the data in the memory after the calculation. If the user would like to execute plot or other data manipulation commands, the save-run-time-result bool should be “True”. However, the user should notice the memory usage when scanning a lot of parameters.

```
Run Time Write Matrix : False
```

Whether to write the results when calculation.

```
Run Time Write Matrix : False  
Run Time Plot : False  
Run Time Contour Plot : False
```

Whether to write the results when calculation.

Plot Setting

```
Plot Region Bool:
  EML : True (EMLz+ add EMLz- )
  EML(EMLz+) : False (at the position just above emitting position)
  EML(EMLz-) : False (at the position just below emitting position)
  +Semi-infinite region: True (at the interface of the 1st semi-infinite region)
  -Semi-infinite region: True (at the interface of the last semi-infinite region)
Plot Mode Bool: (should be considered with plot region bool)
  Total (M1+M2+M3+M4) : True
  +z (M1+M2) : False
  -z (M3+M4) : False
  M1 : False
  M2 : False
  M3 : False
  M4 : False
Write Region Bool:
  EML : True (EMLz+ add EMLz- )
  EML(EMLz+) : True (at the position just above emitting position)
  EML(EMLz-) : True (at the position just below emitting position)
  +Semi-infinite region: True (at the interface of the 1st semi-infinite region)
  -Semi-infinite region: True (at the interface of the last semi-infinite region)
Write Mode Bool: (should be considered with plot region bool)
  Total (M1+M2+M3+M4) : True
  +z (M1+M2) : True
  -z (M3+M4) : True
  M1 : True
  M2 : True
  M3 : True
  M4 : True
WriteMatrix Region Bool:
  EML : True (EMLz+ add EMLz- )
  EML(EMLz+) : False (at the position just above emitting position)
  EML(EMLz-) : False (at the position just below emitting position)
  +Semi-infinite region: True (at the interface of the 1st semi-infinite region)
  -Semi-infinite region: True (at the interface of the last semi-infinite region)
WriteMatrix Mode Bool: (should be considered with plot region bool)
  Total (M1+M2+M3+M4) : True
  +z (M1+M2) : False
  -z (M3+M4) : False
  M1 : False
  M2 : False
```

Calculate Power Density

Bottom-emitting OLED

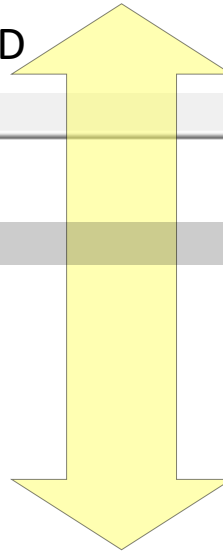
sfile-1.txt		
1	MATERIAL	THICKNESS (nm)
2	air	X
3	Al	120.000000
4	LiF	1.000000
5	B3PYMPM	50
6	CBP	20.000000
7	TAPC	50
8	cito	80
9	glass	X
10		
11		
12		
13		
14	wavelength(nm) : 480.0:10.0:600.0	
15	EML :	
16	Layerno	: 5
17	Position(nm)	: 10.000000
18	ratio	: 1.000000
19	QY	: 1.000000
20	Fluorescence	: cbp_irppy2acac
21	DOF	: cbp_irppy2acac
22		

Thick metal

structureList-PDCmdl1.txt			
1	Name	readfilename	readfilepath
2	#s1	sfile-1.txt	../Example/structure/Convention
	savefilename		
	sfile-1	../Example/PDCm	

Calculate Power Density

Transparent/ Double-emitting OLED



1	MATERIAL	THICKNESS (nm)
2	air	X
3	Al	10:5:20
4	LiF	1.000000
5	B3PYMPM	50
6	CBP	20.000000
7	TAPC	20
8	cito	50
9	glass	X
10		
11		
12		
13	wavelength(nm) :	450:10:650.0
14	EML :	
15	Layerno	: 5
16	Position (nm)	: 10.000000
17	ratio	: 1.000000
18	QY	: 1.000000
19	Fluorescence	: cbp_irppy2acac
20	DOF	: cbp_irppy2acac

Thin metal

1	Name	readfilename	readfilepath	savefilename
2	#Ts1	sfile-1.txt	./Example/structure/Transparent	sfile-1



Calculate Power Density at EML

```
>>> changefigshowBool F
>>> save_run_time_result_Bool F
>>> ReadStructListPath ./Example/structure/Convention
>>> ReadStructListName structureList-PDCmd1.txt
>>> readStructList
Now reading structure list file ./Example/structure/Convention\structureList-PDCmd1.txt


| No./Name | filename    | savefilename | CommandID | Check | readfilepath                                         |
|----------|-------------|--------------|-----------|-------|------------------------------------------------------|
| #s1      | sfile-1.txt | sfile-1      | 0.0       | X     | ./Example/structure/Convention ../../Example/PDCmd-1 |


Structure file reading...
Now reading structure file ./Example/structure/Convention\sfile-1.txt

>>> ReadStructListPath ./Example/structure/Transparent
>>> ReadStructListName structureList-PDCmd1.txt
>>> readStructList
Now reading structure list file ./Example/structure/Transparent\structureList-PDCmd1.txt


| No./Name | filename    | savefilename | CommandID | Check | readfilepath                                         |
|----------|-------------|--------------|-----------|-------|------------------------------------------------------|
| #Ts1     | sfile-1.txt | sfile-1      | 0.0       | X     | ./Example/structure/Transparent ../../Example/PDCmd- |


Structure file reading...
Now reading structure file ./Example/structure/Transparent\sfile-1.txt
```



Calculate Power Density at EML

```
>>> printStructInfo
```

```
*****
```

```
Name: #s1
```

```
[#] Material Thickness(nm)
```

```
-----  
[1] air X  
[2] Al 120.0  
[3] LiF 1.0  
[4] B3PYMPM 50.0  
[5] CBP 20.0  
[6] TAPC 50.0  
[7] cito 80.0  
[8] glass X
```

```
wavelength(nm) : 480.00000:10.00000:600.00000
```

```
EML Fluorescence DOF Position(nm) PLQY Ratio
```

```
-----  
[5] cbp_irppy2acac cbp_irppy2acac 10.0 1.0 1.0
```

```
Device number : 1
```

```
*****
```

```
Name: #Ts1
```

```
[#] Material Thickness(nm)
```

```
-----  
[1] air X  
[2] Al 10.00000:5.00000:20.00000  
[3] LiF 1.0  
[4] B3PYMPM 50.0  
[5] CBP 20.0  
[6] TAPC 20.0  
[7] cito 50.0  
[8] glass X
```

```
wavelength(nm) : 450.00000:10.00000:650.00000
```

```
EML Fluorescence DOF Position(nm) PLQY Ratio
```

```
-----  
[5] cbp_irppy2acac cbp_irppy2acac 10.0 1.0 1.0
```

```
Device number : 3
```

Calculate Power Density at EML

```
*** Unknown syntax: #####  
>>> ktko      0:0.001:2.5  
>>> Thetakt   0  
>>> runtime_write_matrix F  
>>> runtime_plot      T  
>>> runtime_contourplot F
```

Calculate Power Density at EML

```
>>> printPDInfo
Setting file path : ../.. /SETTING/user-1
Setting file name : PDCmd_setting
kt/ko : 0.00000:0.00100:2.50000
theta in kt/ko domain : 0.0
kx/ko : -1.00000:0.10000:1.00000
ky/ko : -1.00000:0.10000:1.00000
z(nm) : -50.00000:1.00000:50.00000
Save Run Time Result : False
Run Time Write Matrix : False
Run Time Plot : True
Run Time Contour Plot : False

Plot Region Bool:
EML : True (EMLz+ add EMLz- )
EML(EMLz+) : False (at the position just above emitting position)
EML(EMLz-) : False (at the position just below emitting position)
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)
-Semi-infinite region: True (at the interface of the last semi-infinite region)
Plot Mode Bool: (should be considered with plot region bool)
Total (M1+M2+M3+M4) : True
+z (M1+M2) : False
-z (M3+M4) : False
M1 : False
M2 : False
M3 : False
M4 : False
Write Region Bool:
EML : True (EMLz+ add EMLz- )
EML(EMLz+) : True (at the position just above emitting position)
EML(EMLz-) : True (at the position just below emitting position)
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)
-Semi-infinite region: True (at the interface of the last semi-infinite region)
```

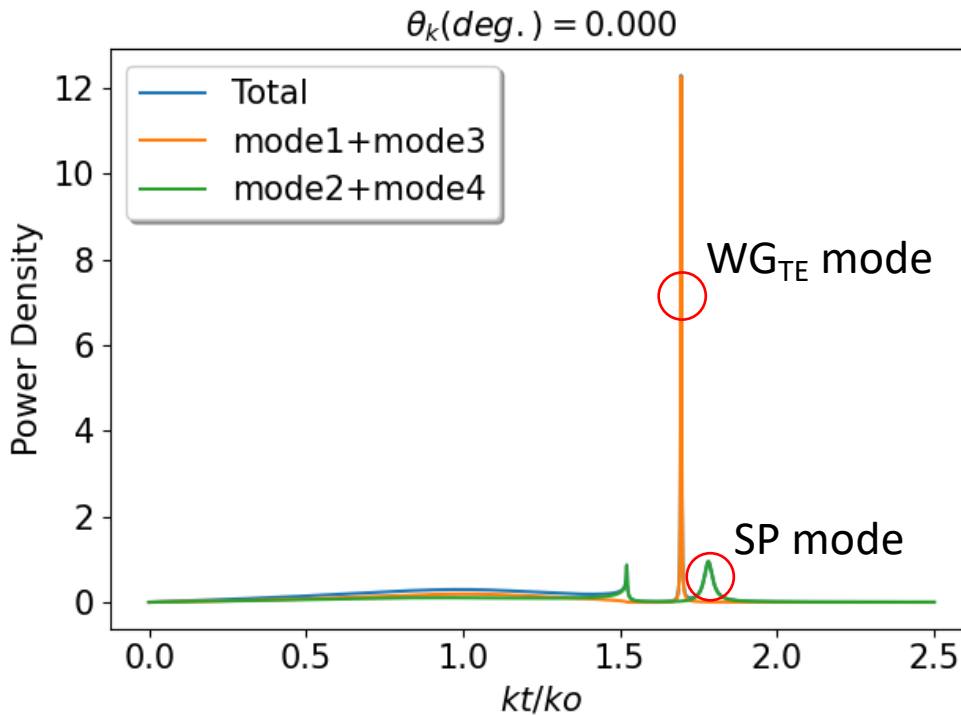
Calculate Power Density at EML

[illegible]

Data at EML

Bottom-emitting OLED

480 nm



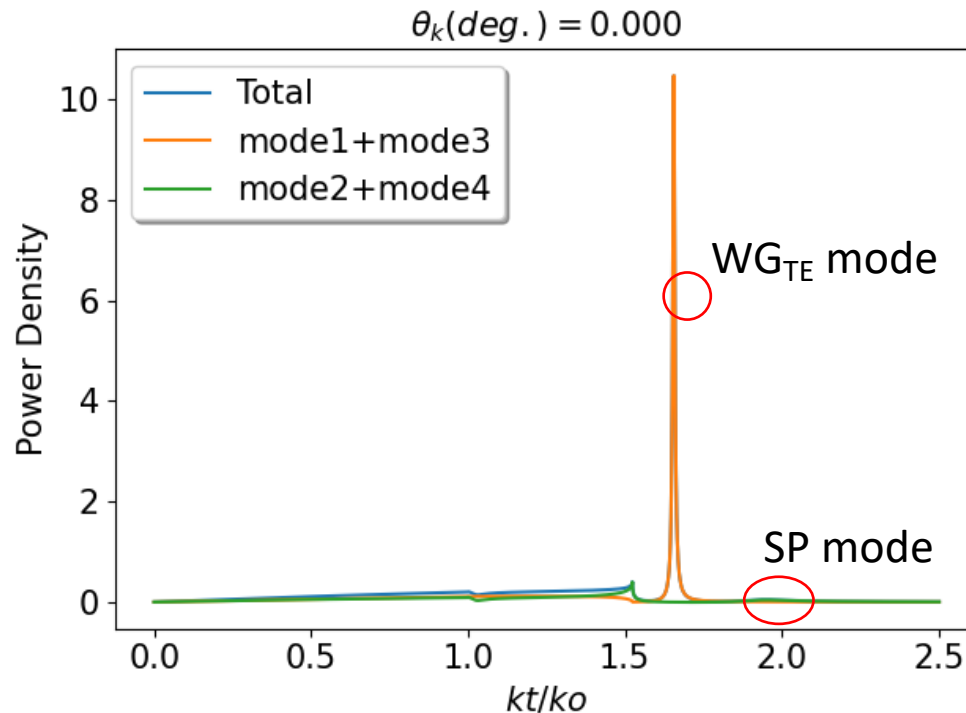
MATERIAL	THICKNESS (nm)
air	X
Al	120.000000
LiF	1.000000
B3PYMPM	50
CBP	20.000000
TAPC	50
cito	80
glass	X

wavelength (nm) : 480.0:10.0:600.0

Data at EML

Transparent/ Double-emitting OLED

480 nm



MATERIAL THICKNESS (nm)	
air	X
Al	10
LiF	1.000000
B3PYMPM	50
CBP	20.000000
TAPC	20
cito	50
glass	X

wavelength (nm) : 480.0:10.0:600.0

Calculate Power Density at EML

```
*** Unknown syntax: #####  
>>> kxko      0:0.01:2.5  
>>> kyko      0:0.01:1  
>>> runtime_write_matrix T  
>>> runtime_plot      F  
>>> runtime_contourplot T
```

Calculate Power Density at EML

```
>>> printPDInfo

Setting file path : ../.. /SETTING/user-1
Setting file name : PDCmd.setting
kt/ko : 0.00000:0.00100:2.50000
Theta in kt/ko domain : 0 0
kx/ko : 0.00000:0.01000:2.50000
ky/ko : 0.00000:0.01000:1.00000
z(nm) : -50.00000:1.00000:50.00000
Save Run Time Result : False
Run Time Write Matrix : True
Run Time Plot : False
Run Time Contour Plot : True

Plot Region Bool:
  EML : True (EMLz+ add EMLz- )
  EML(EMLz+) : False (at the position just above emitting position)
  EML(EMLz-) : False (at the position just below emitting position)
  +Semi-infinite region: True (at the interface of the 1st semi-infinite region)
  -Semi-infinite region: True (at the interface of the last semi-infinite region)
Plot Mode Bool: (should be cosidered with plot region bool)
  Total (M1+M2+M3+M4) : True
  +z (M1+M2) : False
  -z (M3+M4) : False
  M1 : False
  M2 : False
  M3 : False
  M4 : False
Write Region Bool:
  EML : True (EMLz+ add EMLz- )
  EML(EMLz+) : True (at the position just above emitting position)
  EML(EMLz-) : True (at the position just below emitting position)
  +Semi-infinite region: True (at the interface of the 1st semi-infinite region)
  -Semi-infinite region: True (at the interface of the last semi-infinite region)
Write Mode Bool: (should be cosidered with plot region bool)
  Total (M1+M2+M3+M4) : True
  +z (M1+M2) : True
  -z (M3+M4) : True
  M1 : True
  M2 : True
  M3 : True
  M4 : True
```

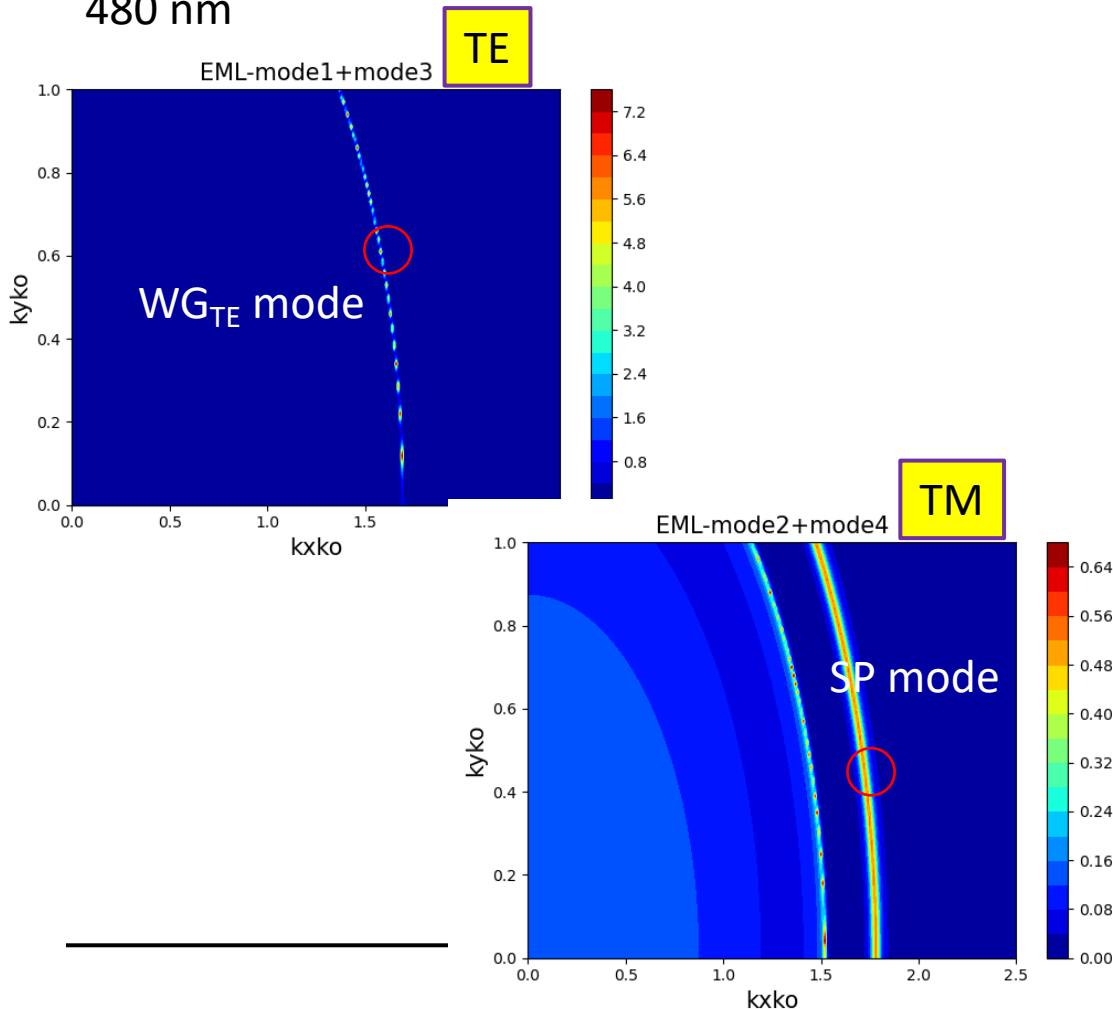

Calculate Power Density at EML

[illegible]

Data at EML

Bottom-emitting OLED

480 nm



MATERIAL	THICKNESS (nm)
air	X
Al	120.000000
LiF	1.000000
B3PYMPM	50
CBP	20.000000
TAPC	50
cito	80
glass	X

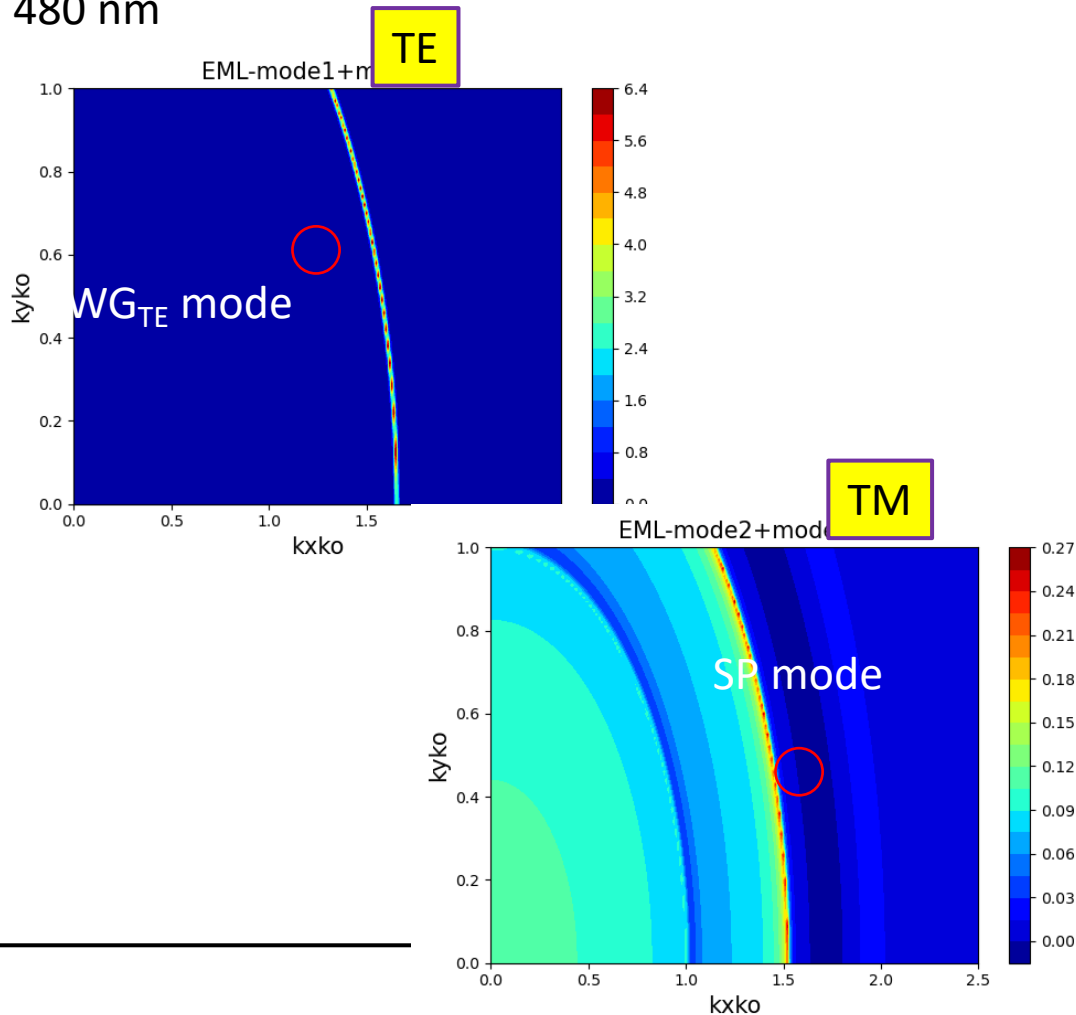
wavelength (nm) : 480.0:10.0:600.0



Data at EML

Transparent/ Double-emitting OLED

480 nm



MATERIAL THICKNESS (nm)		
air		X
Al		10
LiF	1.000000	
B3PYMPM	50	
CBP	20.000000	
TAPC	20	
cito	50	
glass		X

wavelength (nm) : 480.0:10.0:600.0



Calculate Power Density

Bottom-emitting OLED

Diagram of a Bottom-emitting OLED structure. The structure is shown as a stack of layers. A large black 'X' is drawn over the Al layer, indicating it is not the emission layer. A yellow arrow points down from the glass substrate, indicating the emission direction. A yellow box labeled 'Thick metal' points to the Al layer.

Layer	MATERIAL	THICKNESS (nm)
1	air	X
2	Al	120.000000
3	LiF	1.000000
4	B3PYMPM	50
5	CBP	20.000000
6	TAPC	50
7	cito	80
8	glass	X

Parameters:

```

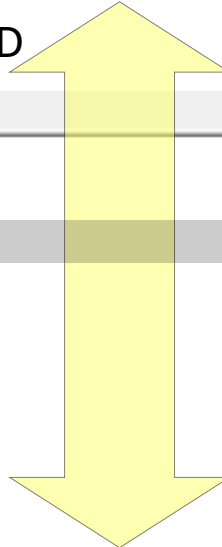
14 wavelength(nm) : 480.0:10.0:600.0
15 EML :
16 Layerno          : 5
17 Position(nm)     : 10.000000
18 ratio            : 1.000000
19 QY               : 1.000000
20 Fluorescence     : cbp_irppy2acac
21 DOF              : cbp_irppy2acac
  
```

Name	readfilename	readfilepath	savefilename
#s1	sfile-1.txt	./Example/structure/Convention	sfile-1

../.../Example/PDzCmd-1/C

Calculate Power Density

Transparent/ Double-emitting OLED



sfile-1.txt	
1	MATERIAL
2	THICKNESS (nm)
2	air
3	Al
4	LiF
5	B3PYMPM
6	CBP
7	TAPC
8	cito
9	glass
10	
11	
12	
13	wavelength(nm) : 450:10:650.0
14	EML :
15	Layerno : 5
16	Position(nm) : 10.000000
17	ratio : 1.000000
18	QY : 1.000000
19	Fluorescence : cbp_irppy2acac
20	DOF : cbp_irppy2acac

Thin metal

ueList-PDeCmd1.txt	
Name	readfilepath
#Ts1	sfile-1.txt
savefilename	sav
../Example/structure/Transparent	sfile-1
../Example/PDzCmd-1/Trans	

Calculate Power Density at z

```
>>> changefigshowBool F
>>> save_run_time_result_Bool F
>>> ReadStructListPath ./Example/structure/Convention
>>> ReadStructListName structureList-PDzCmd1.txt
>>> readStructList

Now reading structure list file ./Example/structure/Convention\structureList-PDzCmd1.txt
  No./Name      filename      savefilename  CommandID  Check      readfilepath
-----
    #s1      sfile-1.txt          sfile-1       0.0       X ./Example/structure/Convention  ../../Example/PDzCmd-1/Conven

Structure file reading...
-----
Now reading structure file ./Example/structure/Convention\sfile-1.txt

>>> ReadStructListPath ./Example/structure/Transparent
>>> ReadStructListName structureList-PDzCmd1.txt
>>> readStructList

Now reading structure list file ./Example/structure/Transparent\structureList-PDzCmd1.txt
  No./Name      filename      savefilename  CommandID  Check      readfilepath
-----
    #Ts1      sfile-1.txt          sfile-1       0.0       X ./Example/structure/Transparent  ../../Example/PDzCmd-1/Trans

Structure file reading...
-----
Now reading structure file ./Example/structure/Transparent\sfile-1.txt
```



Calculate Power Density at z

```

Structure file reading...
-----
Now reading structure file ./Example/structure/Transparent\sfile-1.txt

>>> printStructInfo

*****
Name: #s1
-----
[1]  Material      Thickness(nm)
-----
[1]  air           X
[2]  Al            120.0
[3]  LiF            1.0
[4]  B3PYMPM       50.0
[5]  CBP           20.0
[6]  TAPC          50.0
[7]  cito          80.0
[8]  glass         X

wavelength(nm) : 480.00000:10.00000:600.00000

EML  Fluorescence  DOF          Position(nm)  PLQY      Ratio
-----
[5]  cbp_irppy2acac  cbp_irppy2acac  10.0          1.0          1.0
Device number : 1

*****
Name: #Ts1
-----
[1]  Material      Thickness(nm)
-----
[1]  air           X
[2]  Al            10.00000:5.00000:20.00000
[3]  LiF            1.0
[4]  B3PYMPM       50.0
[5]  CBP           20.0
[6]  TAPC          20.0
[7]  cito          50.0
[8]  glass         X

wavelength(nm) : 450.00000:10.00000:650.00000

EML  Fluorescence  DOF          Position(nm)  PLQY      Ratio
-----
[5]  cbp_irppy2acac  cbp_irppy2acac  10.0          1.0          1.0
Device number : 3
  
```

Calculate Power Density at z

```
Unknown syntax: ~~~~~  
>>> ktko      (0.5,1,1.5)  
>>> Thetakt    0  
>>> z          -50:1:300  
>>> runtime_plot  T
```


Calculate Power Density at z

```
>>> printPDInfo

Setting file path : .././SETTING/user-1
Setting file name : PDCmd.setting
kt/ko : 0.50000:0.50000:1.50000
Theta in kt/ko domain : 0.0
kx/ko : 0.00000:0.01000:2.50000
ky/ko : 0.00000:0.01000:1.00000
z(nm) : -50.00000:1.00000:300.00000
Save Run Time Result : False
Run Time Write Matrix : True
Run Time Plot : True
Run Time Contour Plot : True

Plot Region Bool:
EML : True (EMLz+ add EMLz- )
EML(EMLz+) : False (at the position just above emitting position)
EML(EMLz-) : False (at the position just below emitting position)
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)
-Semi-infinite region: True (at the interface of the last semi-infinite region)
Plot Mode Bool: (should be considered with plot region bool)
Total (M1+M2+M3+M4) : True
+z (M1+M2) : False
-z (M3+M4) : False
M1 : False
M2 : False
M3 : False
M4 : False
Write Region Bool:
EML : True (EMLz+ add EMLz- )
EML(EMLz+) : True (at the position just above emitting position)
EML(EMLz-) : True (at the position just below emitting position)
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)
-Semi-infinite region: True (at the interface of the last semi-infinite region)
Write Mode Bool: (should be considered with plot region bool)
Total (M1+M2+M3+M4) : True
+z (M1+M2) : True
-z (M3+M4) : True
M1 : True
M2 : True
M3 : True
M4 : True
```

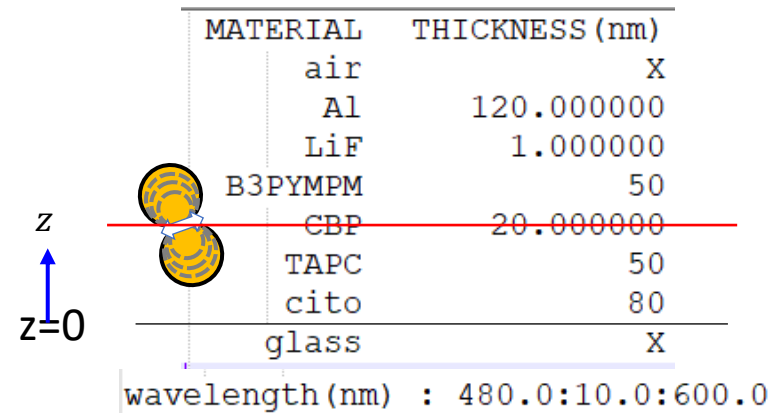
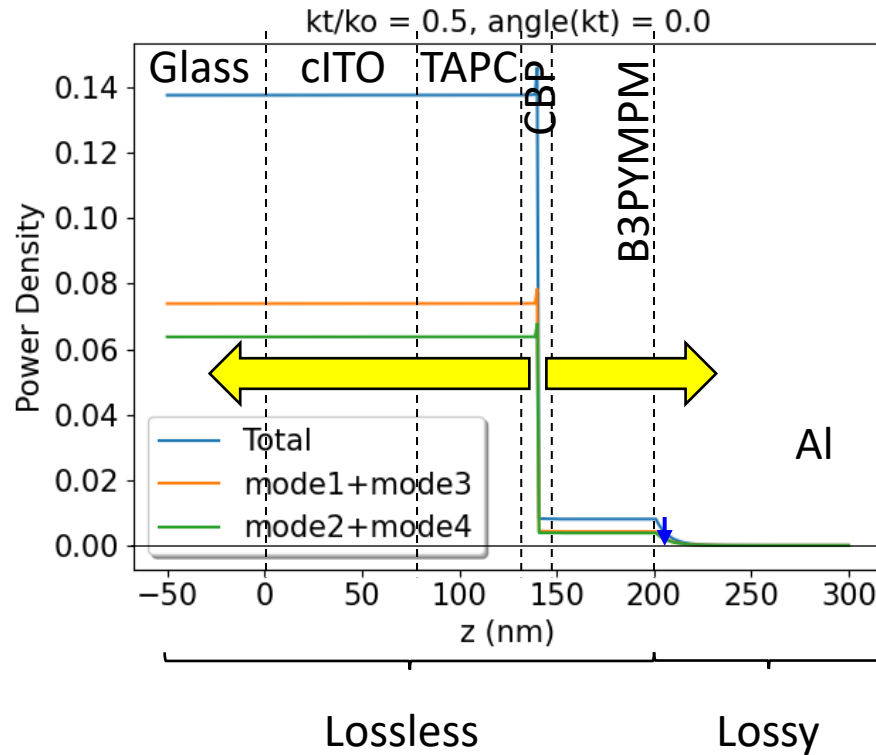
Calculate Power Density at z

[illegible]

Data at EML

Bottom-emitting OLED

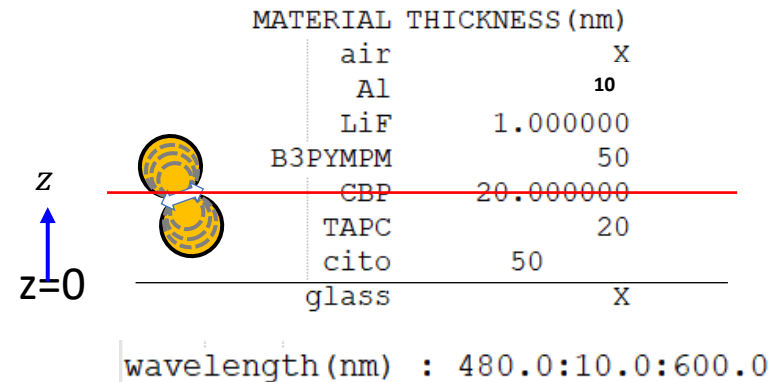
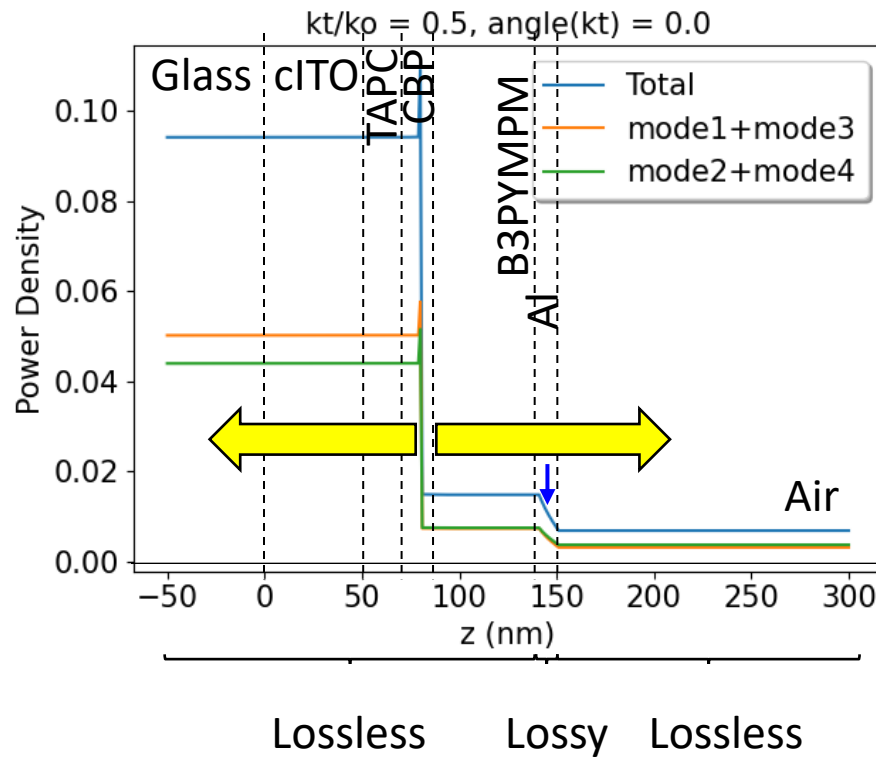
480 nm



Data at EML

Transparent/ Double-emitting OLED

480 nm



Calculate Incoherent Power Density at EML

```
>>> changefigshowBool F
>>> save_run_time_result_Bool F
>>> ReadStructListPath ./Example/structure/Convention
>>> ReadStructListName structureList-PDIncohCmd1.txt
>>> readStructList
Now reading structure list file ./Example/structure/Convention\structureList-PDIncohCmd1.txt
  No./Name      filename      savefilename CommandID Check      readfilepath
-----
    #s1  sfile-1-air.txt          sfile-1      0.0      X ./Example/structure/Convention ../../Example/PDIncohCmd-1/Convention
Structure file reading...
Now reading structure file ./Example/structure/Convention\sfile-1-air.txt

>>> ReadStructListPath ./Example/structure/Transparent
>>> ReadStructListName structureList-PDIncohCmd1.txt
>>> readStructList
Now reading structure list file ./Example/structure/Transparent\structureList-PDIncohCmd1.txt
  No./Name      filename      savefilename CommandID Check      readfilepath
-----
    #Tsa1  sfile-1-air.txt          sfile-1-air    0.0      X ./Example/structure/Transparent ../../Example/PDIncohCmd-1/Transparent
Structure file reading...
Now reading structure file ./Example/structure/Transparent\sfile-1-air.txt
```

Calculate Incoherent Power Density at EML

sfile-1-air.txt	
1	MATERIAL
2	THICKNESS (nm)
2	air
3	X
3	Al
4	120.000000
4	LiF
5	1.000000
5	B3PYMPM
6	50
6	CBP
7	20.000000
7	TAPC
8	50
8	cito
9	80
9	glass
10	X
10	air
11	X
12	
13	
14	wavelength(nm) : 480.0:10.0:600.0
15	EML :
16	Layer no : 5
17	Position(nm) : 10.000000
18	ratio : 1.000000
19	QY : 1.000000
20	Fluorescence : cbp_irppy2acac
21	DOF : cbp_irppy2acac

Calculate Incoherent Power Density at EML

sfile-1-air.txt	
sfile-1-air.txt	
1	MATERIAL THICKNESS (nm)
2	air X
3	Al 10:5:20
4	LiF 1.000000
5	B3PYMPM 50
6	CBP 20.000000
7	TAPC 20
8	cito 50
9	glass X
10	air X
11	
12	
13	wavelength(nm) : 450:10:650.0
14	EML :
15	Layerno : 5
16	Position(nm) : 10.000000
17	ratio : 1.000000
18	QY : 1.000000
19	Fluorescence : cbp_irppy2acac
20	DOF : cbp_irppy2acac

Calculate Incoherent Power Density at EML

```
>>> printStructInfo
```

```
*****
```

```
Name: #s1
```

```
[#] Material Thickness(nm)
```

[1]	air	X
[2]	Al	120.0
[3]	LiF	1.0
[4]	B3PYMPM	50.0
[5]	CBP	20.0
[6]	TAPC	50.0
[7]	cito	80.0
[8]	glass	X
[9]	air	X

```
wavelength(nm) : 480.00000:10.00000:600.00000
```

EML	Fluorescence	DOF	Position(nm)	PLQY	Ratio
-----	--------------	-----	--------------	------	-------

[5]	cbp_irppy2acac	cbp_irppy2acac	10.0	1.0	1.0
-----	----------------	----------------	------	-----	-----

```
Device number : 1
```

```
*****
```

```
Name: #Ts1
```

```
[#] Material Thickness(nm)
```

[1]	air	X
[2]	Al	10.00000:5.00000:20.00000
[3]	LiF	1.0
[4]	B3PYMPM	50.0
[5]	CBP	20.0
[6]	TAPC	20.0
[7]	cito	50.0
[8]	glass	X
[9]	air	X

```
wavelength(nm) : 450.00000:10.00000:650.00000
```

EML	Fluorescence	DOF	Position(nm)	PLQY	Ratio
-----	--------------	-----	--------------	------	-------

[5]	cbp_irppy2acac	cbp_irppy2acac	10.0	1.0	1.0
-----	----------------	----------------	------	-----	-----

```
Device number : 3
```


Calculate Incoherent Power Density at EML

```
*** Unknown syntax: #####  
>>> ktko      0:0.001:2.5  
>>> Thetakt    0  
  
>>> runtime_write_matrix F  
  
>>> runtime_plot      T  
  
>>> runtime_contourplot F  
  
>>> printPDInfo  
  
Setting file path : ../SETTING/user-1  
Setting file name : PDCmd.setting  
kt/ko : 0.00000:0.00100:2.50000  
Theta in kt/ko domain : 0.0  
kx/ko : 0.00000:0.01000:2.50000  
ky/ko : 0.00000:0.01000:1.00000  
z(nm) : -50.00000:1.00000:300.00000  
Save Run Time Result : False  
Run Time Write Matrix : False  
Run Time Plot : True  
Run Time Contour Plot : False  
  
Plot Region Bool:  
EML : True (EMLz+ add EMLz- )  
EML(EMLz+) : False (at the position just above emitting position)  
EML(EMLz-) : False (at the position just below emitting position)  
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)  
-Semi-infinite region: True (at the interface of the last semi-infinite region)  
Plot Mode Bool: (should be considered with plot region bool)  
Total (M1+M2+M3+M4) : True  
+z (M1+M2) : False  
-z (M3+M4) : False  
M1 : False  
M2 : False  
M3 : False  
M4 : False  
  
Write Region Bool:  
EML : True (EMLz+ add EMLz- )  
EML(EMLz+) : True (at the position just above emitting position)  
EML(EMLz-) : True (at the position just below emitting position)  
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)  
-Semi-infinite region: True (at the interface of the last semi-infinite region)  
Write Mode Bool: (should be considered with plot region bool)  
Total (M1+M2+M3+M4) : True  
+z (M1+M2) : True  
-z (M3+M4) : True  
M1 : True  
M2 : True
```

Calculate Incoherent Power Density at EML

```

Semi-Infinite Region: True (at the interface of the last semi-infinite region)
Write Mode Bool: (should be considered with plot region bool)
  Total (M1+M2+M3+M4) : True
  +z (M1+M2) : True
  -z (M3+M4) : True
  M1 : True
  M2 : True
  M3 : True
  M4 : True
WriteMatrix Region Bool:
  EML : True (EMLz+ add EMLz- )
  EML(EMLz+) : False (at the position just above emitting position)
  EML(EMLz-) : False (at the position just below emitting position)
  +Semi-infinite region: True (at the interface of the 1st semi-infinite region)
  -Semi-infinite region: True (at the interface of the last semi-infinite region)
WriteMatrix Mode Bool: (should be considered with plot region bool)
  Total (M1+M2+M3+M4) : True
  +z (M1+M2) : False
  -z (M3+M4) : False
  M1 : False
  M2 : False
  M3 : False
  M4 : False
Figure Show Bool : False

```

[illegible]

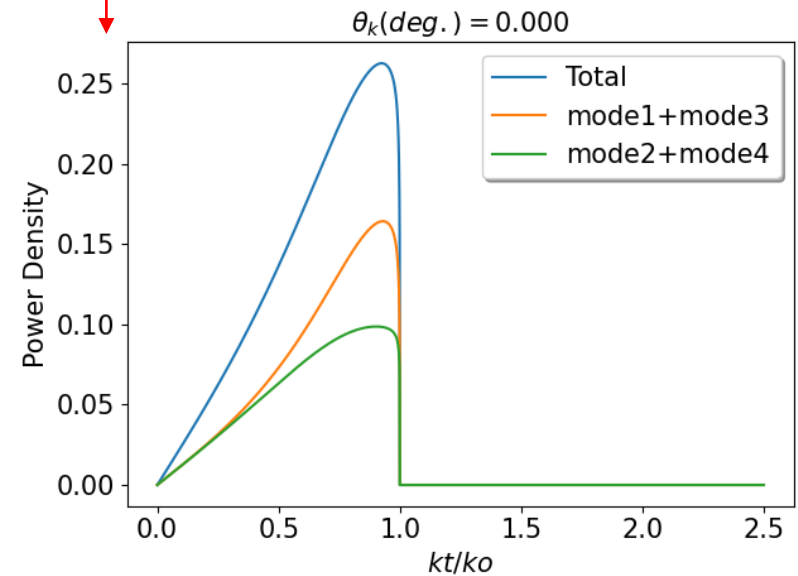
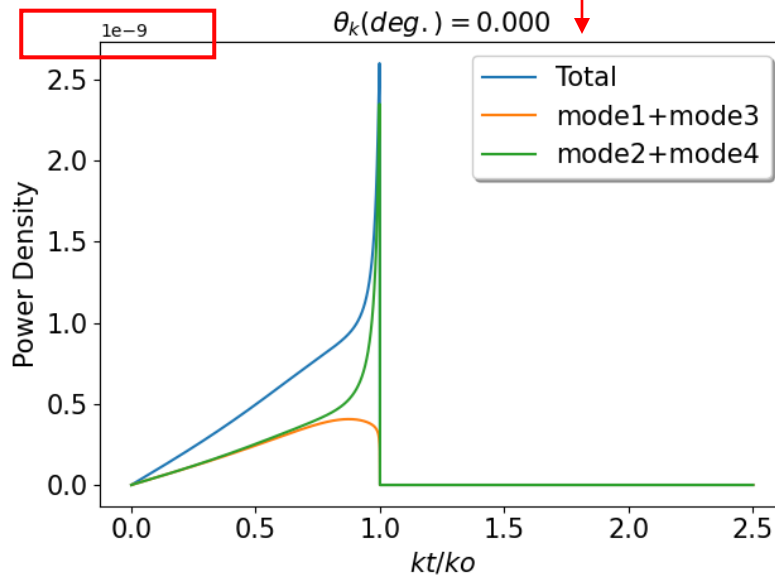
Data at EML

Bottom-emitting OLED

480 nm

MATERIAL	THICKNESS (nm)
air	X
Al	120.000000
LiF	1.000000
B3PYMPM	50
CBP	20.000000
TAPC	50
cito	80
glass	X
air	X

wavelength (nm) : 480.0:10.0:600.0

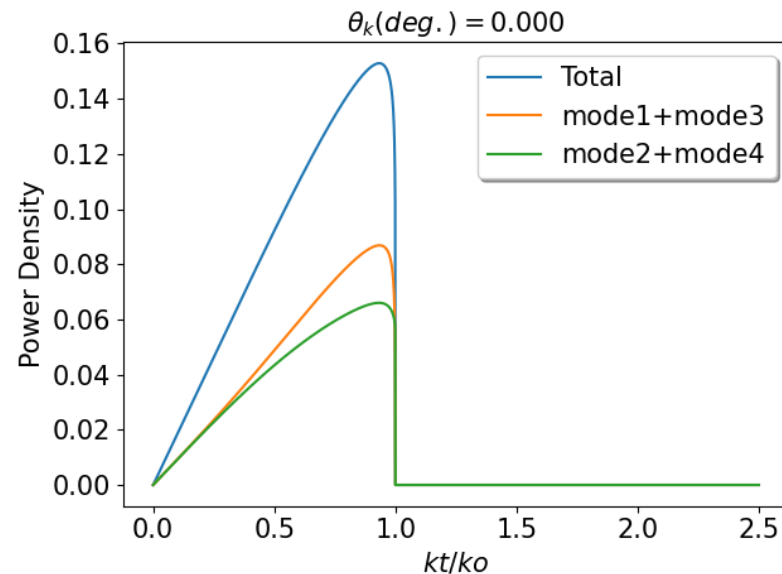
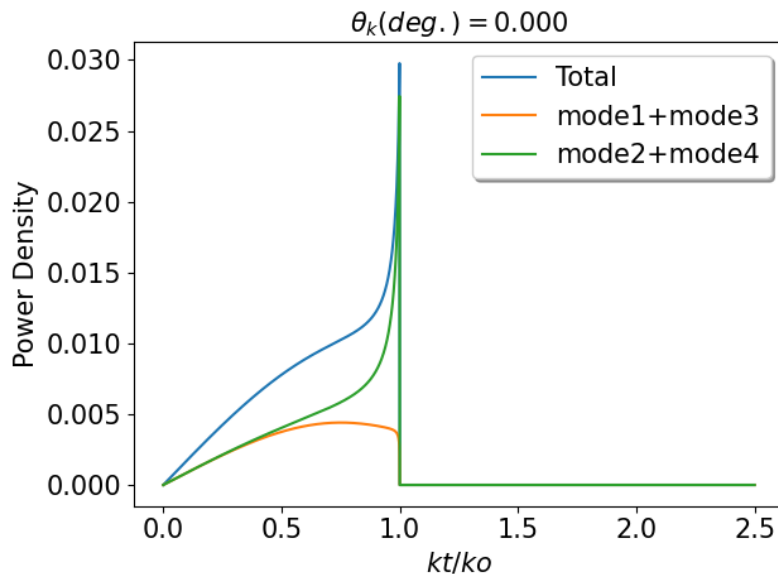


Data at EML

Transparent/ Double-emitting OLED

480 nm

MATERIAL	THICKNESS (nm)
air	X
Al	10:5:20
LiF	1.000000
B3PYMPM	50
CBP	20.000000
TAPC	20
cito	50
glass	X
air	X
wavelength (nm) : 480.0:10.0:600.0	



Calculate Incoherent Power Density at EML

```
>>> kxko      0:0.01:2.5
>>> kyko      0:0.01:1
>>> runtime_write_matrix T
>>> runtime_plot      F
>>> runtime_contourplot T
>>> printPDInfo

Setting file path : .././SETTING/user-1
Setting file name : PDCmd.setting
kt/ko : 0.00000:0.00100:2.50000
Theta in kt/ko domain : 0.0
kx/ko : 0.00000:0.01000:2.50000
ky/ko : 0.00000:0.01000:1.00000
z(nm) : -50.00000:1.00000:300.00000
Save Run Time Result : False
Run Time Write Matrix : True
Run Time Plot : False
Run Time Contour Plot : True

Plot Region Bool:
EML : True (EMLz+ add EMLz- )
EML(EMLz+) : False (at the position just above emitting position)
EML(EMLz-) : False (at the position just below emitting position)
+Semi-infinite region: True (at the interface of the 1st semi-infinite region)
-Semi-infinite region: True (at the interface of the last semi-infinite region)
Plot Mode Bool: (should be considered with plot region bool)
Total (M1+M2+M3+M4) : True
+z (M1+M2) : False
-z (M3+M4) : False
M1 : False
M2 : False
M3 : False
M4 : False
```

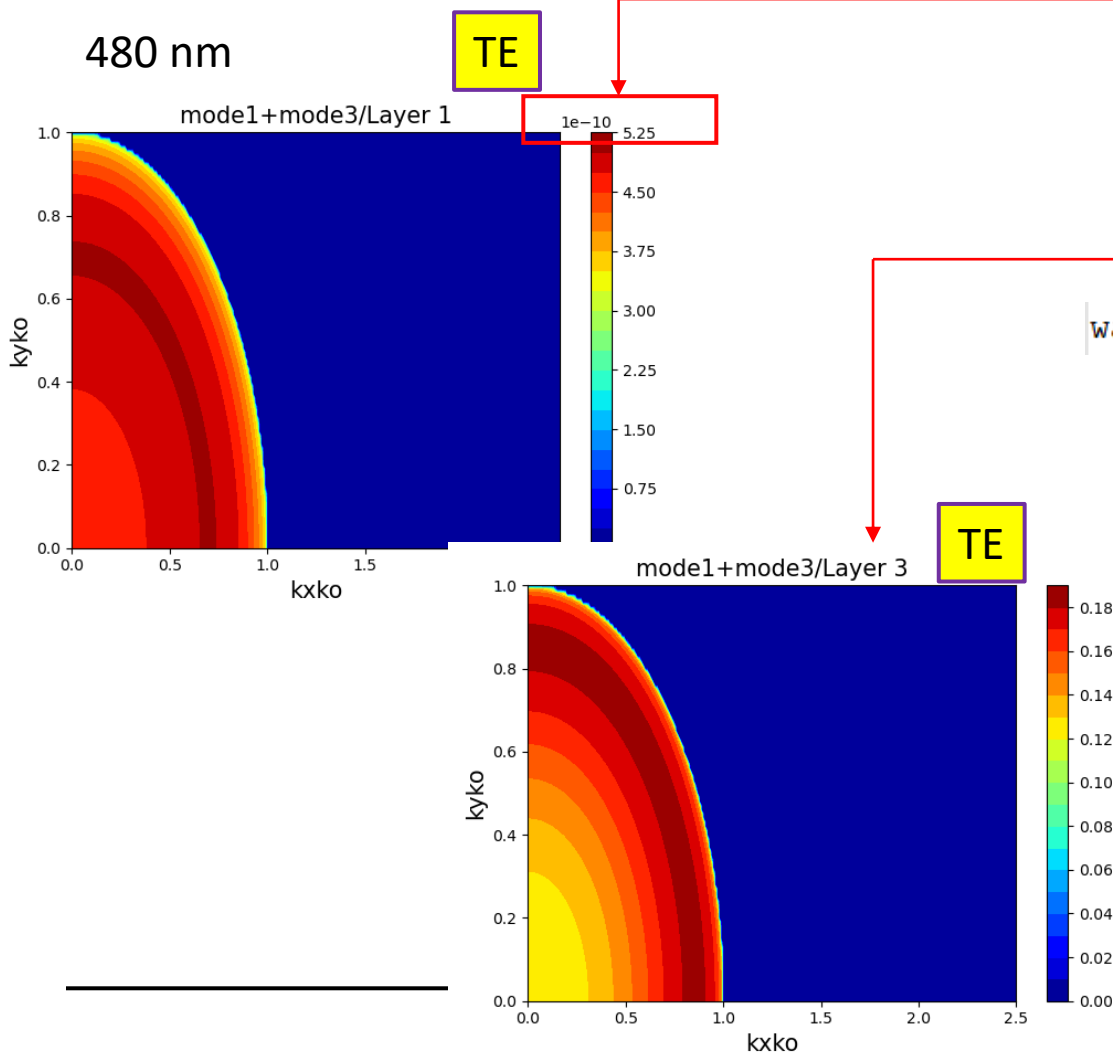
Calculate Incoherent Power Density at EML

[illegible]

Data at EML

Bottom-emitting OLED

480 nm



MATERIAL	THICKNESS (nm)
air	X
Al	120.000000
LiF	1.000000
B3PYMPM	50
CBP	20.000000
TAPC	50
cito	80
glass	X
air	X

wavelength (nm) : 480.0:10.0:600.0

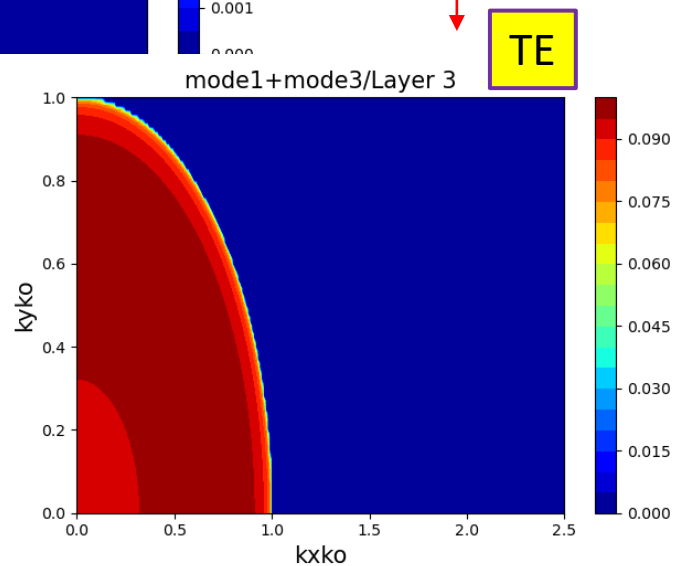
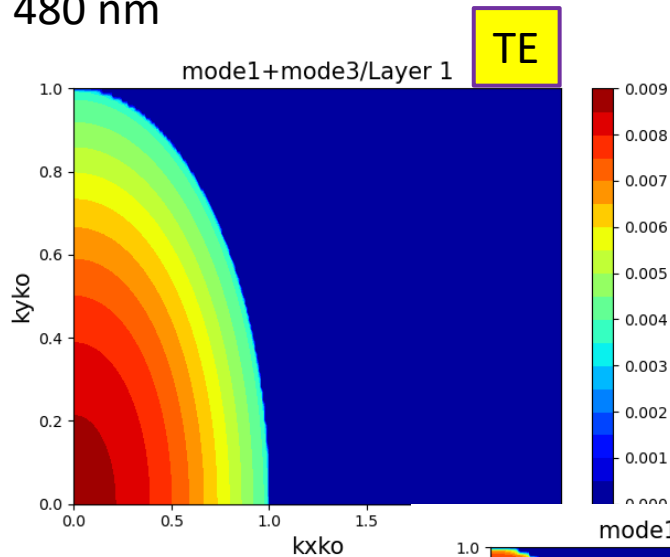
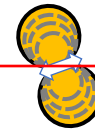


Data at EML

Transparent/ Double-emitting OLED

480 nm

MATERIAL	THICKNESS (nm)
air	X
Al	10:5:20
LiF	1.000000
B3PYMPM	50
CBP	20.000000
TAPC	20
cito	50
glass	X
air	X



wavelength (nm) : 480.0:10.0:600.0



Calculate Device Distribution

Exit the material manager system.

*** The material manager system would be automatically saved into the user's setting directory.

```
>>> exit
```

```
-----  
End running GOODLAB anisotropic simulator ver1.0 Tue Apr 14 10:56:58 2020  
Elapsed time : 0 day(s)/ 1 hr(s)/ 6 min(s)/ 3.5559067726195254 sec(s)  
-----
```

