

## Chapter 14

# DOSE RESPONSE OF DOSIMETRIC MATERIALS: MODELS

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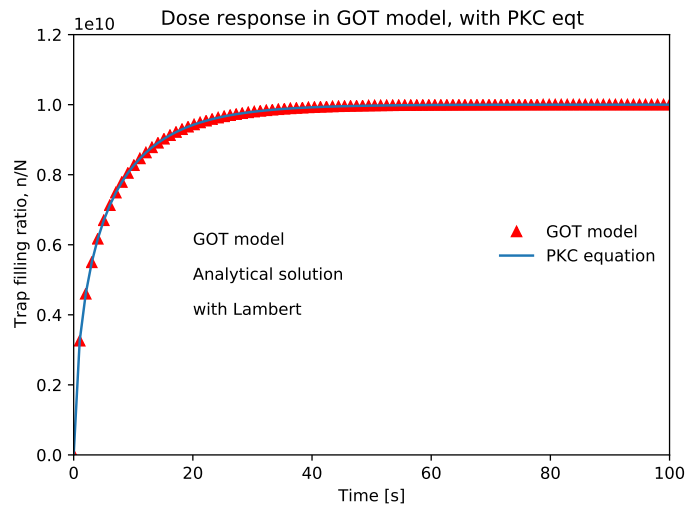
**Code 14.1: Irradiation: OTOR, Lambert analytical solution**

```
# GOT MODEL- code for IRRADIATION Lambert solution
import numpy as np
from scipy.integrate import odeint
import matplotlib.pyplot as plt
import warnings
warnings.filterwarnings("ignore")
from scipy.special import lambertw
# Numerical parameters for GOT irradiation model.
N, R, X, n0 = \
1E10, .1, 1e10, 0
t = np.linspace(0, 100, 100)
# The GOT model irradiation differential equation.
def deriv(y, t):
    n = y
    dndt = (N-n) * X*R / ((N - n) * R + n )
    return dndt
y0 = n0
ret = odeint(deriv, y0, t)
n = ret.flatten()
plt.plot(t, n, 'r^', lw=2, label='GOT model');
ret = odeint(deriv, y0, t)
#analytical n(t)
Dc=N/R
u=N*(1+np.real(lambertw((R-1)*np.exp((R-1)-\
(t*X/np.abs(Dc))))/(1-R)))
plt.plot(t, u,label='PKC equation');
```

```

leg = plt.legend(loc='right')
leg.get_frame().set_linewidth(0.0)
plt.ylabel('Trap filling ratio, n/N');
plt.ylim(0,1.2e10);
plt.xlim(0,100);
plt.title('Dose response in GOT model, with PKC eqt');
plt.text(20,.6e10,'GOT model');
plt.text(20,.5e10,'Analytical solution');
plt.text(20,.4e10,'with Lambert');
plt.xlabel('Time [s]');
plt.tight_layout()
plt.show()

```



**Fig. 14.1:** Dose response in the OTOR model, using the Lambert analytical solution PKC Eq.(??). The analytical solution (solid line) is compared with the numerical solution of the differential Eq.(??) shown as triangles.

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#### Code 14.2: Plots of the PKC equation

```

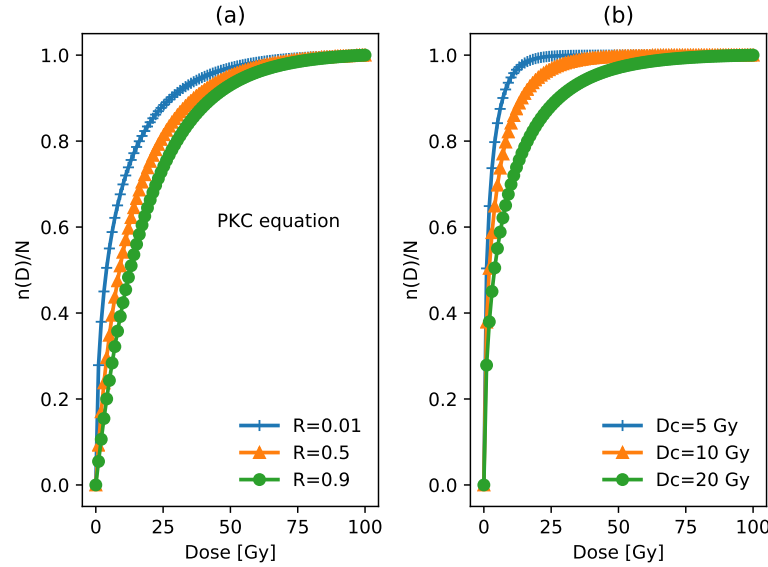
#plot dose response using PKC equation for different R, Dc
import numpy as np

```

```

import matplotlib.pyplot as plt
from scipy.special import lambertw
import warnings
warnings.filterwarnings("ignore")
def PKC(R,Dc):
    u=(1+np.real(lambertw((R-1)*np.exp((R-1)-\
    (D/np.abs(Dc)))))/(1-R))
    plt.plot(D,u/max(u),syms[j-1], linewidth=2,
    label=labls[j-1]);
Dc=20
D = np.linspace(0, 100, 100)
Rs=[.01,.5,.9]
labls=["R="+str(x) for x in Rs]
syms=['+-','^-','o-']
plt.subplot(1,2,1);
for j in range(1,4):
    PKC(Rs[j-1],Dc)
leg = plt.legend()
leg.get_frame().set_linewidth(0.0)
plt.xlabel('Dose [Gy]');
plt.text(45,.6,'PKC equation');
plt.ylabel('n(D)/N');
plt.title('(a)');
plt.subplot(1,2,2);
Dcs=[5,10,20]
labls=['Dc='+str(x)+' Gy' for x in Dcs]
for j in range(1,4):
    PKC(0.01,Dcs[j-1])
leg = plt.legend()
leg.get_frame().set_linewidth(0.0)
plt.xlabel('Dose [Gy]');
plt.ylabel('n(D)/N');
plt.title('(b)');
plt.tight_layout()
plt.show()

```



**Fig. 14.2:** (a) Plot of the PKC Eq.(??) for 3 values of the re trapping ratio  $R = 0.01, 0.5, 0.9$  and a fixed characteristic dose  $D_c = 20$  Gy. As the value of  $R$  increases, the dose response curve takes longer to reach saturation, due to re trapping (b) Plot of the PKC equation for 3 values of  $D_c = 5, 10, 20$  Gy and a fixed  $R = 0.01$ . As the value of  $D_c$  increases, the dose response curve takes longer to reach saturation.

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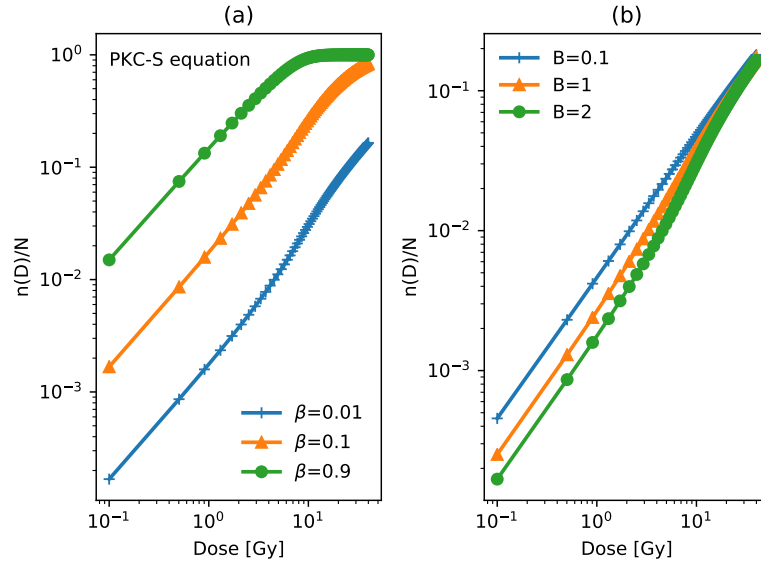
### Code 14.3: Plots of the PKC-S equation

```
#plot dose response PKC-S equation for different beta, B
import numpy as np
import matplotlib.pyplot as plt
from scipy.special import lambertw
import warnings
warnings.filterwarnings("ignore")
def PKCS(B,beta):
    u=np.real(lambertw(B*np.exp(np.abs(B)-(D/Dc))))/B
    u=(1-(u**beta))
    plt.plot(D,u,syms[j-1], linewidth=2,
    label=labls[j-1]);
```

```

        plt.yscale("log");
        plt.xscale("log");
B, Dc= 2, 2
D = np.linspace(.1, 40, 100)
betas=[.01,.1,.9]
labls=[r'$\beta$'+str(x) for x in betas]
syms=['+-','^-', 'o-']
plt.subplot(1,2,1);
for j in range(1,4):
    PKCS(B,betas[j-1])
leg = plt.legend()
leg.get_frame().set_linewidth(0.0)
plt.xlabel('Dose [Gy]');
plt.text(.1,.8,'PKC-S equation')
# plt.text(45,.6,'PKC equation');
plt.ylabel('n(D)/N');
plt.title('(a)');
plt.subplot(1,2,2);
Bs=[.1,1,2]
labls=['B'+str(x) for x in Bs]
for j in range(1,4):
    PKCS(Bs[j-1],.01)
leg = plt.legend()
leg.get_frame().set_linewidth(0.0)
plt.xlabel('Dose [Gy]');
plt.ylabel('n(D)/N');
plt.title('(b)');
plt.tight_layout()
plt.show()

```



**Fig. 14.3:** (a) Plot of the PKC-S Eq.(??) for 3 values of the constant  $\beta = 0.01, 0.1, 0.9$  and a fixed constant  $B = 2$ . As the value of  $\beta$  increases, the amount of superlinearity decreases. (b) Plot of the PKC-S equation for  $B = 0.1, 1, 2$  and a fixed  $\beta = 0.01$ . As the value of  $B$  increases, the amount of superlinearity increases. For more details see Pagonis et al. [1]

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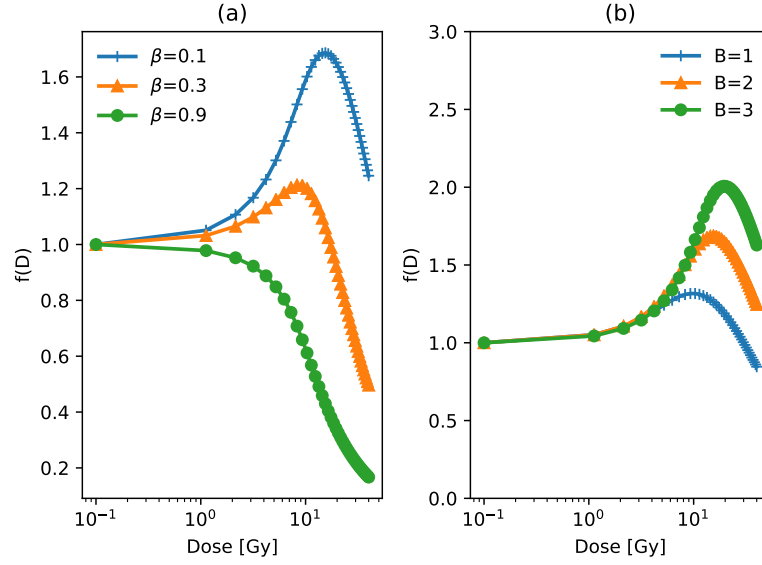
#### Code 14.4: Plots of the supralinearity index $f(D)$

```
#plot coefficient f(D) for different beta, B
import numpy as np
import matplotlib.pyplot as plt
from scipy.special import lambertw
import warnings
warnings.filterwarnings("ignore")
def PKCS(B,beta):
    u=np.real(lambertw(B*np.exp(np.abs(B)-(D/Dc))))/B
    u=(1-(u**beta))/D
    u1=np.real(lambertw(B*np.exp(np.abs(B)-(.1/Dc))))/B
    u1=(1-(u1**beta))/.1
    plt.plot(D,u/u1,symbs[j-1], linewidth=2,
```

```

        label=labls[j-1]);
        plt.xscale("log");
B, Dc= 2, 2
D = np.linspace(.1, 40, 40)
betas=[.1,.3,.9]
labls=[r'$\beta$'+str(x) for x in betas]
syms=['+-','^-','o-']
plt.subplot(1,2,1);
for j in range(1,4):
    PKCS(B,betas[j-1])
leg = plt.legend()
leg.get_frame().set_linewidth(0.0)
plt.xlabel('Dose [Gy]');
plt.ylabel('f(D)');
plt.title('(a)');
plt.subplot(1,2,2);
Bs=[1,2,3]
labls=['B'+str(x) for x in Bs]
for j in range(1,4):
    PKCS(Bs[j-1],.1)
leg = plt.legend()
leg.get_frame().set_linewidth(0.0)
plt.ylim(0,3);
plt.xlabel('Dose [Gy]');
plt.ylabel('f(D)');
plt.title('(b)');
plt.tight_layout()
plt.show()

```



**Fig. 14.4:** (a) Plot of Eq.(??) for 3 values of the constant  $\beta = 0.1, 0.3, 0.9$  and a fixed constant  $B = 2$ . As the value of  $\beta$  increases, the value of  $f(D)$  and the amount of superlinearity decreases. (b) Plot of Eq.(??) for  $B = 1, 2, 3$  and a fixed  $\beta = 0.1$ . As the value of  $B$  increases, the value of  $f(D)$  and the amount of superlinearity increases. For more details see Pagonis et al. [1]



## References

1. V. Pagonis, G. Kitis, R. Chen, Superlinearity revisited: A new analytical equation for the dose response of defects in solids, using the Lambert W function, *Journal of Luminescence* 227 (2020) 117553. doi:<https://doi.org/10.1016/j.jlumin.2020.117553>. URL <http://www.sciencedirect.com/science/article/pii/S0022231320310449>