

```

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% INITIALIZE %%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
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%%% https://research.tuni.fi/metaplasmonics/
%%% +358 50 447 8330

close all;
clc;
clear;

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% DATA IMPORTATION %%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
load dipole_dipole_normalized.mat; %% Save the normalized decay rate and the dipole-dipole interactions
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% Here we used the above name

r = lum.x0; %% interatomic distance
gamma = lum.y1; %% decay rate gamma_12
gparam_norm = lum.y0; %% dipole-dipole interactions -- g12

```

## Selection of some section of concurrence to plot as a function of time

```

indt1 = 200; %% INDEX PLOT FOR TIME CONCURRENCE NOT WITH POSITION
indt2 = 413;

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% DASHBOARD %%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
time_value = 20; %% time -- normalized time with gamma11 ==1
t = linspace(0, time_value, 413);
gg = gamma(ceil(length(gamma)/2)); %% self interactions g11

```

## concurrence section

```

B = exp(-(gg+gamma).*t'); %% evolution of sub and superradiant density matrix
C = exp(-(gg-gamma).*t');
D = (B-C).^2;
E = 4*exp(-(2.*gg.*t)).*sin(2.*gparam_norm'.*t).^2;
Conc = 0.5*sqrt(D + E. '); %% concurrence formulations

gamma_0 = gamma(ceil(length(gamma)/2));
tt = linspace (0,2,431);
tau = tt;

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% PLOTTING SECTION %%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%

```

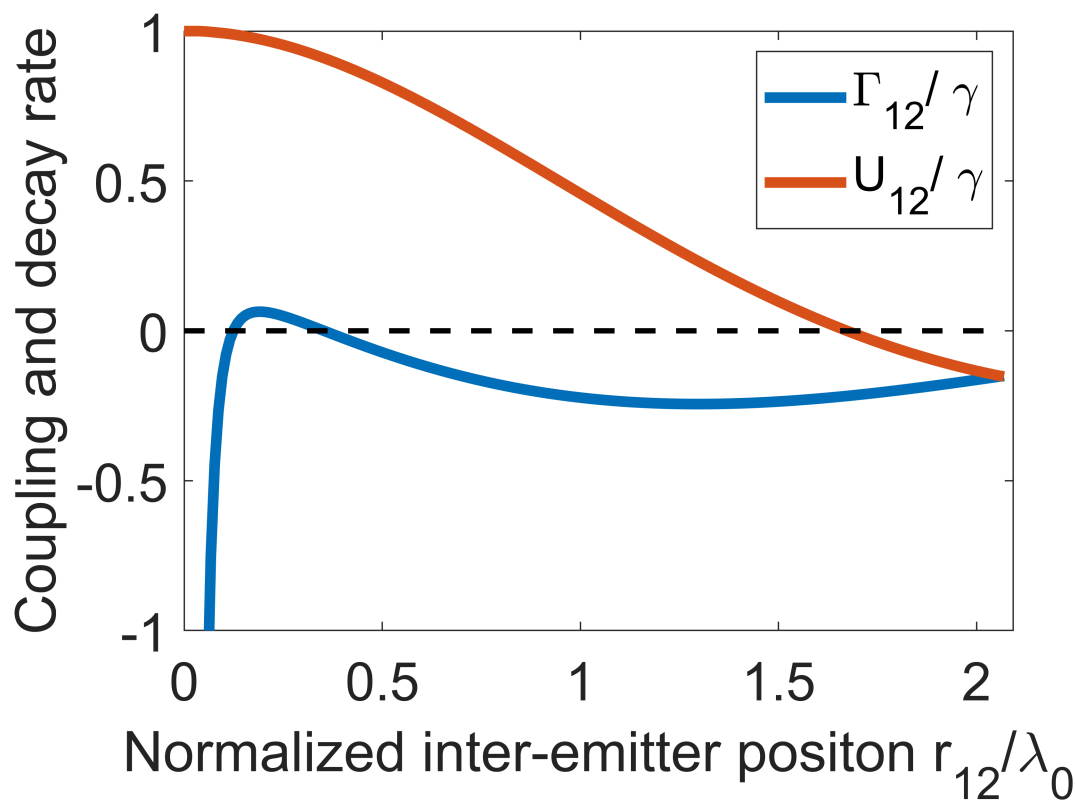
```

%% POWER PLOT %%%%%%%%%%%
%
rho_ee = exp(-tau');
rho_aa = exp(-(gamma_0-gamma).*tau');
rho_ss =exp(-(gamma_0+gamma).*tau');

P = 2*gamma_0*rho_ee + (gamma_0 + gamma).*rho_ss + (gamma_0 - gamma).*rho_aa;

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% PLOTTING SECTION %%%%%%%%%%%
fig2 = figure('color', 'w');
ax1 = axes;
plot(lum.x0, lum.y0,lum.x1, lum.y1, 'Linewidth', 4)
hold on
plot(lum.x0, 0*lum.y0, 'k--', 'Linewidth', 2);
set(ax1, 'XLim', [0 2.09351])
set(ax1, 'YLim', [-1 1])
set(gca, 'FontSize', 20);
xlabel('Normalized inter-emitter positon  $r_{12}/\lambda_0$ ', 'FontSize',21);
ylabel('Coupling and decay rate', 'FontSize', 21);
legend('\Gamma_{12}/ \gamma', 'U_{12}/ \gamma')

```



```

% Create figure
figure1 = figure('color', 'w');
% Create axes

```

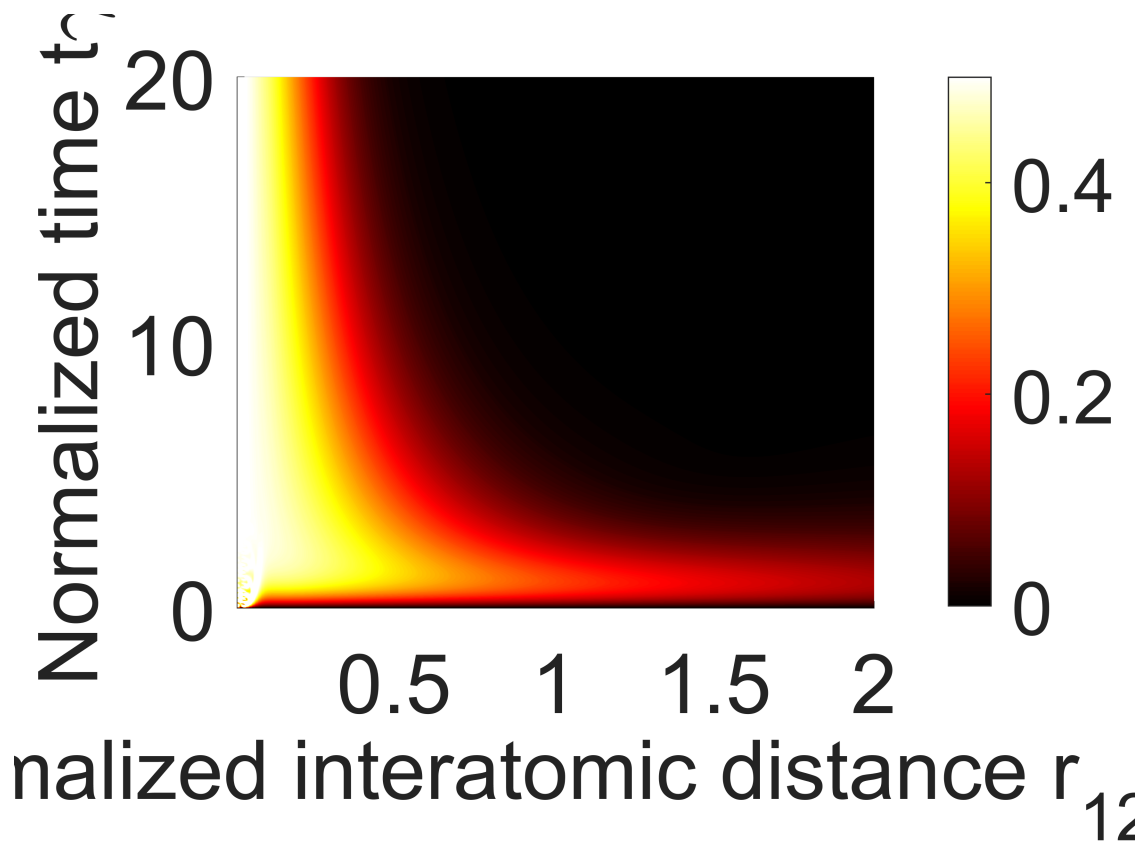
```

axes1 = axes('Parent',figure1,'FontSize',31);
view(axes1,[90 -90]);
grid(axes1,'on');
hold(axes1,'all');
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% PLOTTING SECTION %%%%%%%%%%
[x, y]=meshgrid(t,r);
surf(x,y,Conc.', 'Parent',axes1,'LineStyle','none',...
      'FaceColor','interp');
% Create colorbar
colorbar('peer',axes1);
colormap('hot');
caxis([0 .5])
ylim([0.01 2])

set(gca,'FontSize', 31);
%xlabel('Normalized interatomic distance  $r_{12}/\lambda_0$ ');
%ylabel('Normalized time  $t/\gamma$ ');
xlabel('Normalized time  $t/\gamma$ ','VerticalAlignment','bottom','Rotation',90,...
      'HorizontalAlignment','center',...
      'FontSize',31);

ylabel('Normalized interatomic distance  $r_{12}/\lambda_0$ ','VerticalAlignment','cap',...
      'HorizontalAlignment','center',...
      'FontSize',31);

```



```

%title('Concurrence (C)');

```

# pOWER PLOT

Create figure

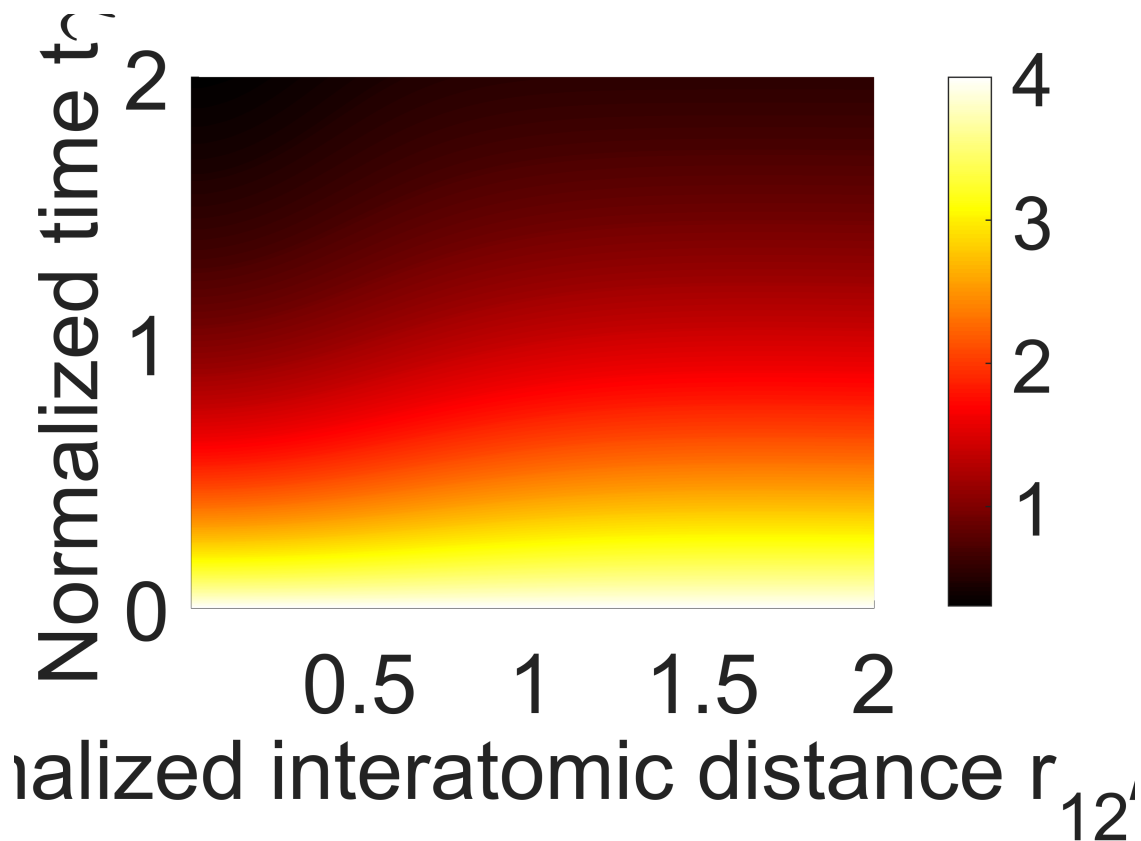
```
figure1 = figure('color', 'w');

% Create axes
axes1 = axes('Parent',figure1,'FontSize',31);
view(axes1,[90 -90]);
grid(axes1,'on');
hold(axes1,'all');
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%% PLOTTING SECTION %%%%%%%%%%%%%%
[x, y]=meshgrid(tt,r);
surf(x,y,P.','Parent',axes1,'LineStyle','none',...
     'FaceColor','interp');

% Create colorbar
colorbar('peer',axes1);
colormap('hot');
%caxis([0.7 1.2])
ylim([0.01 2])
xlim([0, .5]);

xlabel('Normalized time  $t \backslash \gamma$ ','VerticalAlignment','bottom','Rotation',90,...
     'HorizontalAlignment','center',...
     'FontSize',31);

ylabel('Normalized interatomic distance  $r_{12} / \lambda_{0}$ ','VerticalAlignment','cap',...
     'HorizontalAlignment','center',...
     'FontSize',31);
```



```
%title('Power enhancement')
```

```
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
```

```
fig4 = figure('color', 'w');
```

```
ax1 = axes;
```

```
plot(lum.x0, Conc(indt1, :), 'Linewidth', 4)
```

```
hold on
```

```
plot(lum.x0, Conc(indt2, :), 'Linewidth', 4)
```

```
%plot(lum.x0, 0*lum.y0, 'k--', 'Linewidth', 2);
```

```
xlim([0.01 2])
```

```
set(gca, 'FontSize', 20);
```

```
xlabel('Normalized inter-emitter position  $r_{12}/\lambda_0$ ', 'FontSize', 21);
```

```
ylabel('C(\infty)', 'FontSize', 21);
```

```
legend(strcat('t = ', num2str(ceil(t(indt1))), 's'), strcat('t = ', num2str(ceil(t(indt2))), 's'))
```

