Untitled11

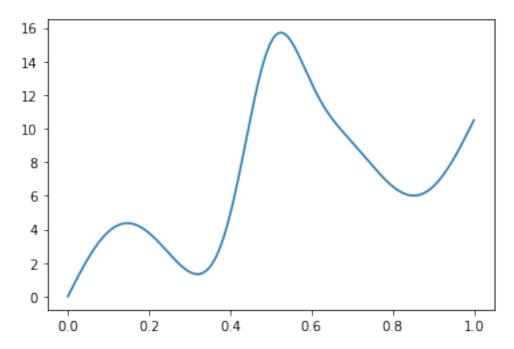
September 29, 2025

1 Une première utilisation de l'intégrale de Lebesgue

```
[15]: import numpy as np
     import matplotlib.pyplot as plt
     # Paramètres
     a, b, c = 0.5, 10, 3
     Left, Right = 0, 1
     N = 100000 # nombre de subdivisions pour Riemann
              # nombre de niveaux pour Lebesque
     # Fonction à intégrer
     def f(x):
         return a*x**2 + b*x + c*np.sin(4*np.pi*x) + 10*np.exp(-100*(x-0.5)**2)
     # Intégrale de Riemann (somme de rectangles)
     # -----
     x_riemann = np.linspace(Left, Right, N)
     dx = (Right - Left)/N
     integral_riemann = np.sum(f(x_riemann) * dx)
     print("Intégrale de Riemann :", integral riemann)
     # Intégrale de Lebesgue (approximation)
     # -----
     # On approxime l'intégrale comme somme sur les niveaux de f
     f_min, f_max = np.min(f(x_riemann)), np.max(f(x_riemann))
     levels = np.linspace(f_min, f_max, M)
     dy = (f_max - f_min)/M
     integral_lebesgue = 0
     for y in levels:
         # mes(\{x \mid f(x) \ge y\}) approximée par fraction de points
         measure = np.sum(f(x riemann) >= y) * dx
         integral_lebesgue += measure * dy
```

```
print("Intégrale de Lebesgue :", integral_lebesgue)
plt.plot(x_riemann,f(x_riemann))
plt.show()
```

Intégrale de Riemann : 6.939103626372618 Intégrale de Lebesgue : 6.9400172353712755



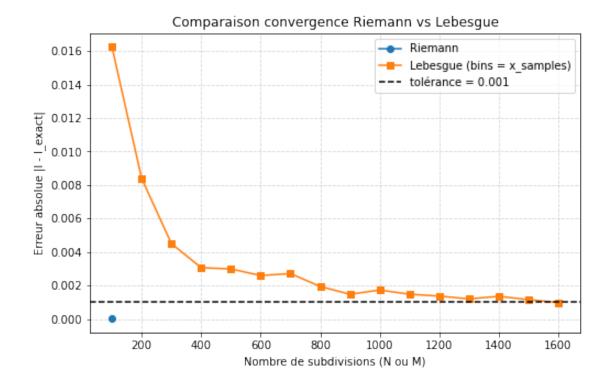
1.1 Raffinement de maillage pour obtenir une erreur de 1e-5

```
I_exact, _ = quad(lambda x: f(x, a, b, c), Left, Right)
print(f"Valeur de référence (quad) {I_exact:.8f}\n")
# MÉTHODE DE RIEMANN (point milieu)
# -----
def riemann_integral(N):
   dx = (Right - Left) / N
   x_vals = Left + dx * (0.5 + np.arange(N))
   return np.sum(f(x_vals, a, b, c)) * dx
N = 100 \# départ
step = 100
errors_riemann = []
Ns_riemann = []
while True:
   I_riem = riemann_integral(N)
   err = abs(I_riem - I_exact)
   errors_riemann.append(err)
   Ns_riemann.append(N)
   print(f"Riemann : N = {N < 5d}) I = {I_riem : .8f} erreur = {err : .2e}")
   if err < tol:</pre>
      N_riem_min = N
      break
   N += step
print(f"\n Riemann : erreur < {tol} atteinte pour N = {N_riem_min} points⊔
 \Rightarrow (erreur = {err:.2e})\n")
# -----
# MÉTHODE DE LEBESGUE (bins = x_samples)
# -----
def lebesgue_integral(M):
   xs = np.linspace(Left, Right, M)
   fx = f(xs, a, b, c)
   f_min, f_max = fx.min(), fx.max()
   edges = np.linspace(f_min, f_max, M + 1) # même nombre de bins
   counts, _ = np.histogram(fx, bins=edges)
   meas = (Right - Left) * counts / len(xs)
   y_lower = edges[:-1]
   y_upper = edges[1:]
   lower = np.sum(y_lower * meas)
   upper = np.sum(y_upper * meas)
   return 0.5 * (lower + upper)
```

```
M = 100
errors_lebesgue = []
Ms_lebesgue = []
while True:
    I_leb = lebesgue_integral(M)
    err = abs(I_leb - I_exact)
    errors_lebesgue.append(err)
    Ms_lebesgue.append(M)
    print(f"Lebesgue : M = {M:<5d} I = {I_leb:.8f} erreur = {err:.2e}")</pre>
    if err < tol:
        M_{leb_min} = M
        break
    M += step
print(f"\n Lebesgue : erreur < {tol} atteinte pour M = {M_leb_min} bins⊔
 \Rightarrow(erreur = {err:.2e})\n")
# GRAPHIQUE COMPARATIF
# -----
plt.figure(figsize=(8,5))
plt.plot(Ns_riemann, errors_riemann, 'o-', label='Riemann')
plt.plot(Ms_lebesgue, errors_lebesgue, 's-', label='Lebesgue (bins =_u
 →x_samples)')
plt.axhline(tol, color='k', linestyle='--', label=f'tolérance = {tol}')
plt.xlabel("Nombre de subdivisions (N ou M)")
plt.ylabel("Erreur absolue |I - I_exact|")
plt.title("Comparaison convergence Riemann vs Lebesgue")
plt.legend()
plt.grid(True, ls="--", alpha=0.5)
plt.show()
Valeur de référence (quad) 6.93912052
Riemann : N = 100
                    I = 6.93911635 erreur = 4.17e-06
 Riemann : erreur < 0.001 atteinte pour N = 100 points (erreur = 4.17e-06)
Lebesgue : M = 100
                     I = 6.92286274 erreur = 1.63e-02
Lebesgue : M = 200
                     I = 6.93073202 erreur = 8.39e-03
Lebesgue : M = 300
                    I = 6.93462358 erreur = 4.50e-03
Lebesgue : M = 400
                    I = 6.93605487 erreur = 3.07e-03
Lebesgue : M = 500
                    I = 6.93613518 erreur = 2.99e-03
                    I = 6.93652614 erreur = 2.59e-03
Lebesgue : M = 600
Lebesgue : M = 700 I = 6.93640871 erreur = 2.71e-03
Lebesgue : M = 800
                    I = 6.93718080 erreur = 1.94e-03
```

```
Lebesgue : M = 900
                     I = 6.93764665 erreur = 1.47e-03
Lebesgue : M = 1000
                     I = 6.93739360 erreur = 1.73e-03
Lebesgue : M = 1100
                     I = 6.93764390 erreur = 1.48e-03
Lebesgue : M = 1200
                     I = 6.93775502 erreur = 1.37e-03
Lebesgue : M = 1300
                     I = 6.93792320 erreur = 1.20e-03
Lebesgue : M = 1400
                     I = 6.93776808 erreur = 1.35e-03
Lebesgue : M = 1500
                     I = 6.93796689 erreur = 1.15e-03
Lebesgue : M = 1600
                     I = 6.93814087 erreur = 9.80e-04
```

Lebesgue : erreur < 0.001 atteinte pour M = 1600 bins (erreur = 9.80e-04)



1.2 Test de l'erreur via une suite de Cauchy

```
a, b, c = 0.5, 10.0, 3.0
tol = 1e-3
# Valeur de référence (juste pour affichage)
I_exact, _ = quad(lambda x: f(x, a, b, c), Left, Right)
print(f"Valeur de référence (quad) {I_exact:.10f}\n")
# Méthode de Lebesque
# -----
def lebesgue_integral(M, M_bins=32):
   xs = np.linspace(Left, Right, M)
    fx = f(xs, a, b, c)
    f_min, f_max = fx.min(), fx.max()
    edges = np.linspace(f_min, f_max, M_bins + 1)
    counts, _ = np.histogram(fx, bins=edges)
    meas = (Right - Left) * counts / len(xs)
    y_lower = edges[:-1]
    y_upper = edges[1:]
    lower = np.sum(y_lower * meas)
    upper = np.sum(y_upper * meas)
    return 0.5 * (lower + upper)
# Boucle avec critère de Cauchy
M = 100
             # maillage initial
M bins = 100 # nombre de bins initial
I_prev = None # valeur précédente
k = 0
while True:
   k += 1
    I_curr = lebesgue_integral(M, M_bins)
    if I_prev is not None:
        diff = abs(I_curr - I_prev)
        print(f"[it=\{k\}] M = \{M: < 6d\} I = \{I\_curr: .10f\} |\Delta I| = \{diff: .2e\}"\}
        if diff < tol:</pre>
            print("\n Critère de Cauchy atteint !")
    else:
        print(f"[it={k}] M = {M:<6d} I = {I_curr:.10f} (première itération)")</pre>
    I_prev = I_curr
    M += 100
    M_bins += 100
```

```
# Résumé
# -----
err = abs(I_curr - I_exact)
print("\n--- Résumé ---")
print(f"Intégrale Lebesgue {I_curr:.10f}")
print(f"Erreur absolue (par rapport à quad) = {err:.2e}")
print(f"Nombre de points M = {M}")
print(f"Nombre de bins M_bins = {M_bins}")
print(f"Itérations = {k}")
Valeur de référence (quad) 6.9391205176
[it=1] M = 100
                 I = 6.9228627416 (première itération)
[it=2] M = 200
                 I = 6.9307320224 |\Delta I| = 7.87e-03
[it=3] M = 300 I = 6.9346235796 |\Delta I| = 3.89e-03
[it=4] M = 400
                 I = 6.9360548702 |\Delta I| = 1.43e-03
[it=5] M = 500
                 I = 6.9361351802 |\Delta I| = 8.03e-05
 Critère de Cauchy atteint!
--- Résumé ---
Intégrale Lebesgue 6.9361351802
Erreur absolue (par rapport à quad) = 2.99e-03
Nombre de points M = 500
```

1.3 Controle de la métrique et critère d'arret de Cauchy

Nombre de bins M_bins = 500

Itérations = 5

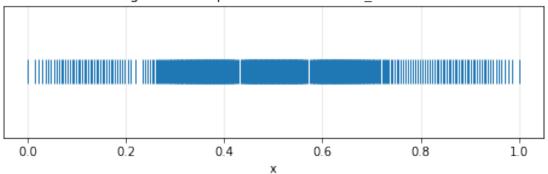
```
# Paramètres
# -----
Left, Right = 0.0, 1.0
a, b, c = 0.5, 10.0, 3.0
tol = 1e-3
                        # critère de Cauchy I_{k+1}-I_k
eps metric = 1e-4
hmin, hmax = 1e-4, 0.2
# (Info) Valeur de référence pour vérifier à la fin
I_exact, _ = quad(lambda x: f(x, a, b, c), Left, Right)
print(f"Référence quad (info) {I_exact:.8f}\n") # devrait être 6.93912052
# Métrique et maillage adapté
# -----
def build curvature mesh (M_cells, eps=eps_metric, hmin=hmin, hmax=hmax,__
 →refN=5000):
    .....
   Construit un maillage adapté \{x_j\}_{j=0...M_cells} (M_cells cellules, _
 \hookrightarrow M cells+1 noeuds)
    en équidistribuant l'arclength dans la métrique sqrt(lambda(x)),
   avec lambda(x) \sim |f''(x)|/eps bornée par [1/hmax^2, 1/hmin^2].
   xr = np.linspace(Left, Right, refN)
   lam = np.abs(f2(xr, a, c)) / eps
   lam = np.clip(lam, 1.0/(hmax**2), 1.0/(hmin**2))
   rho = np.sqrt(lam)
   s = np.cumsum(0.5*(rho[1:]+rho[:-1]) * np.diff(xr))
   s = np.insert(s, 0, 0.0)
   S = s[-1]
   targets = np.linspace(0.0, S, M cells+1)
   x_nodes = np.interp(targets, s, xr)
   return x_nodes # longueur M_cells+1
# Intégrale de Lebesgue pondérée (maillage non uniforme)
# -----
def lebesgue_weighted(nodes, nbins):
   11 11 11
   nodes: array des noeuds (taille M_cells+1)
   nbins: nombre de bins sur l'axe des valeurs (on prend nbins = M_cells)
   - On évalue f au milieu de chaque cellule
   - Poids = longueur de cellule dx
   - Histogramme pondéré par dx -> mesure(E_i)
```

```
HHHH
   x_nodes = nodes
   dx = np.diff(x_nodes)
                                               # poids (mesure locale)
   x_mid = 0.5*(x_nodes[:-1] + x_nodes[1:]) # milieux
   y_mid = f(x_mid, a, b, c)
   ymin, ymax = y_mid.min(), y_mid.max()
   pad = 1e-12 * (abs(ymin)+abs(ymax)+1.0)
   edges = np.linspace(ymin - pad, ymax + pad, nbins+1)
   \# mesure(E_i) = somme des dx des cellules dont y_mid tombe dans le bin
   meas, _ = np.histogram(y_mid, bins=edges, weights=dx)
   y_lower = edges[:-1]
   y_upper = edges[1:]
   lower = np.sum(y_lower * meas)
   upper = np.sum(y_upper * meas)
   I = 0.5*(lower + upper)
   gap = upper - lower # utile si tu veux aussi contrôler upper-lower
   return I, gap
# -----
# Boucle Cauchy + maillage métrique (M = M_bins)
# -----
M \text{ cells} = 100
I_prev = None
k = 0
nodes_final = None
I_curr = None
while True:
   k += 1
   nbins = M_cells # contrainte M = M_bins (ici nbins = nb cellules)
   nodes = build_curvature_mesh(M_cells)
   I_curr, gap = lebesgue_weighted(nodes, nbins)
   if I_prev is not None:
        diff = abs(I_curr - I_prev)
        print(f''[it=\{k\}] M=\{M\_cells:<5d\} I=\{I\_curr:.8f\} |\Delta I|=\{diff:.2e\} |_{U} 

¬(gap={gap:.2e})")
        if diff < tol:</pre>
           print("\n Critère de Cauchy atteint.")
           nodes_final = nodes
            break
```

```
else:
        print(f"[it={k}] M={M_cells:<5d} I={I_curr:.8f} (première itération)")</pre>
    I_prev = I_curr
    M_cells += 100 # raffinement progressif
# Résumé & visualisation finale
# -----
err_abs = abs(I_curr - I_exact)
print("\n--- Résumé ---")
print(f"I (Lebesgue pondéré) {I_curr:.8f}")
print(f"Erreur vs quad (info) = {err_abs:.2e}")
print(f"Maillage final: M_cells = {M_cells} (bins = {M_cells})")
print(f"Itérations = {k}")
plt.figure(figsize=(8,2))
plt.plot(nodes_final, np.zeros_like(nodes_final), '|', markersize=20)
plt.title(f"Maillage final adapté à la courbure (M_cells = {M_cells})")
plt.xlabel("x")
plt.yticks([])
plt.grid(True, alpha=0.3)
plt.show()
Référence quad (info) 6.93912052
[it=1] M=100
               I=6.93793484 (première itération)
[it=2] M=200
               I=6.94178210 |\Delta I|=3.85e-03 (gap=7.54e-02)
[it=3] M=300 I=6.93745650 |ΔI|=4.33e-03 (gap=5.08e-02)
               I=6.94005895 |ΔI|=2.60e-03 (gap=3.83e-02)
[it=4] M=400
[it=5] M=500 I=6.93952161 |ΔI|=5.37e-04 (gap=3.07e-02)
 Critère de Cauchy atteint.
--- Résumé ---
I (Lebesgue pondéré) 6.93952161
Erreur vs quad (info) = 4.01e-04
Maillage final: M_cells = 500 (bins = 500)
Itérations = 5
```

Maillage final adapté à la courbure (M_cells = 500)



2 Approximation de Pi via des calculs d'intégrales

```
[5]: import math
     import random
     # Fonction à intégrer
     f = lambda x: 4/(1+x**2)
     a, b = 0, 1
     # Valeur exacte via erf
     I_exact = math.pi
     # Méthode Riemann
     def riemann_integral(f, a, b, N):
         dx = (b - a)/N
         total = 0.0
         for i in range(N):
             xi = a + i*dx
             total += f(xi)*dx
         return total
     # Méthode Lebesgue
     def lebesgue_integral(f, a, b, Nx, Ny):
         x_vals = [a + i*(b-a)/Nx for i in range(Nx)]
         f_vals = [f(x) for x in x_vals]
         f_{max} = max(f_{vals})
         dy = f_max/Ny
         total = 0.0
         for j in range(Ny):
             yj = j*dy
             measure = 0.0
             for x in x_vals:
                 if f(x) > yj:
```

```
measure += (b-a)/Nx
       total += measure*dy
   return total
# Méthode Monte-Carlo
def monte_carlo_integral(f, a, b, N):
   total = 0.0
   for _ in range(N):
       x = a + (b-a)*random()
       total += f(x)
   return (b-a)*total/N
# Paramètres
N_riemann = 10
Nx, Ny = 10, 10
N_mc = 10
# Calcul des intégrales
I_riemann = riemann_integral(f, a, b, N_riemann)
I_lebesgue = lebesgue_integral(f, a, b, Nx, Ny)
I_mc = monte_carlo_integral(f, a, b, N_mc)
# Affichage du tableau comparatif
print("{:<15} {:<20} ".format("Méthode", "Intégrale approx.", "Erreur

¬relative"))
print("-"*55)
print("{:<15} {:<20.8f} ".format("Riemann", I_riemann,__
 ⇒abs(I_riemann-I_exact)/I_exact))
print("{:<15} {:<20.8f} ".format("Lebesgue", I_lebesgue",
 ⇒abs(I_lebesgue-I_exact)/I_exact))
print("{:<15} {:<20.8f} ".format("Monte-Carlo", I_mc, __
 →abs(I_mc-I_exact)/I_exact))
print("{:<15} {:<20.8f}".format("Exacte", I_exact, 0.0))</pre>
```

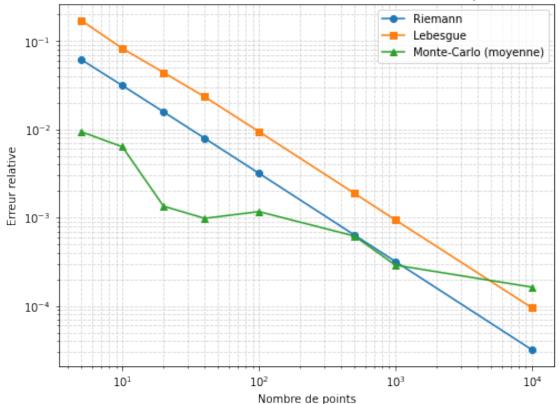
```
Méthode
                    Intégrale approx.
                                         Erreur relative
    Riemann
                    3.23992599
                                         0.03130047
                                         0.08225361
    Lebesgue
                    3.40000000
    Monte-Carlo
                                         0.02591846
                    3.06016741
    Exacte
                    3.14159265
[4]: import math
    import random
    import matplotlib.pyplot as plt
    # Fonction à intégrer
    f = lambda x: 4/(1+x**2)
```

a, b = 0, 1

```
# Valeur exacte via erf
I_exact = math.pi
# Méthode Riemann
def riemann_integral(f, a, b, N):
    dx = (b - a)/N
    total = 0.0
    for i in range(N):
        xi = a + i*dx
        total += f(xi)*dx
    return total
# Méthode Lebesque
def lebesgue_integral(f, a, b, Nx, Ny):
    x_vals = [a + i*(b-a)/Nx for i in range(Nx)]
    f_vals = [f(x) for x in x_vals]
    f_{max} = max(f_{vals})
    dy = f_max/Ny
    total = 0.0
    for j in range(Ny):
        yj = j*dy
        measure = 0.0
        for x in x vals:
            if f(x) > yj:
               measure += (b-a)/Nx
        total += measure*dy
    return total
# Méthode Monte-Carlo moyenne sur M tirages
def monte_carlo_integral(f, a, b, N, M=10):
    total = 0.0
    for _ in range(M):
        s = 0.0
        for _ in range(N):
            x = a + (b-a)*random.random()
            s += f(x)
        total += (b-a)*s/N
    return total/M
# Liste des nombres de points
N_{points} = [5, 10, 20, 40, 100, 500, 1000, 10000]
# Tableaux pour stocker les erreurs relatives
err_riemann = []
err_lebesgue = []
err_mc = []
```

```
for N in N_points:
   I_r = riemann_integral(f, a, b, N)
   I_l = lebesgue_integral(f, a, b, Nx=N, Ny=N)
   I_m = monte_carlo_integral(f, a, b, N, M=50)
   err_riemann.append(abs(I_r - I_exact)/I_exact)
   err_lebesgue.append(abs(I_1 - I_exact)/I_exact)
   err_mc.append(abs(I_m - I_exact)/I_exact)
# Plot des erreurs relatives
plt.figure(figsize=(8,6))
plt.loglog(N_points, err_riemann, 'o-', label='Riemann')
plt.loglog(N_points, err_lebesgue, 's-', label='Lebesgue')
plt.loglog(N_points, err_mc, '^-', label='Monte-Carlo (moyenne)')
plt.xlabel('Nombre de points')
plt.ylabel('Erreur relative')
plt.title('Evolution de l\'erreur relative en fonction du nombre de points')
plt.grid(True, which='both', linestyle='--', alpha=0.5)
plt.legend()
plt.show()
```

Evolution de l'erreur relative en fonction du nombre de points



3 ADRS_multiple_mesh_adapt.py

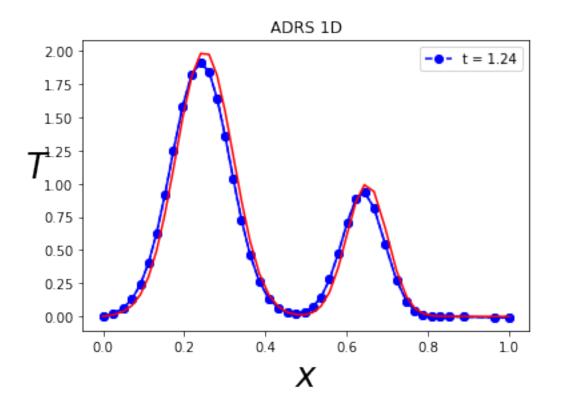
```
[13]: import math
      import numpy as np
      import matplotlib.pyplot as plt
      def adrs_fct(n, x):
      \#u, t = -V u, x + k u, xx - lamda u + f
         u=np.ones(n)
         return u
      def metric_fct(n, u):
      #calcul metric hloc
         hloc=np.ones(n)
         return hloc
      def mesh_fct(n, hloc):
      #calcul metric hloc
         x=np.ones(n)
         return x
      iplot=0
      # PHYSICAL PARAMETERS
      K = 0.01 #Diffusion coefficient
      xmin = 0.0
      xmax = 1.0
      Time = 10. #Integration time
      V=1.
      lamda=1
      #mesh adaptation param
      niter_refinement=30 #niter different calculations
      hmin=0.02
      hmax=0.15
      err=0.03
      # NUMERICAL PARAMETERS
               #Number of grid points : initialization
      NX = 3
      NT = 10000 #Number of time steps max
      ifre=1000000 #plot every ifre time iterations
      eps=0.001
                 #relative convergence ratio
```

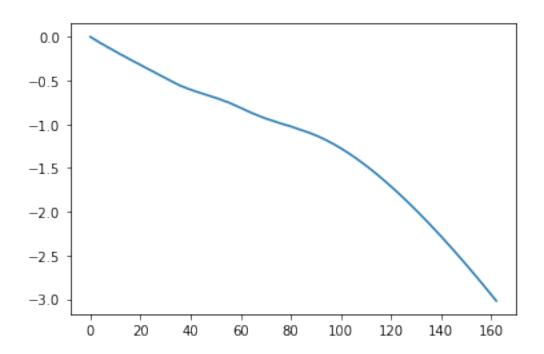
```
errorL2=np.zeros((niter_refinement))
errorH1=np.zeros((niter_refinement))
itertab=np.zeros((niter_refinement))
hloc = np.ones((NX))*hmax
itera=0
NXO=0
while( np.abs(NXO-NX) > 2 and itera<niter_refinement-1):</pre>
    itera+=1
    itertab[itera]=1./NX
    iplot=itera-2
    x = np.linspace(xmin,xmax,NX)
    T = np.zeros((NX))
#mesh adaptation using local metric
    if(itera>0):
        xnew=[]
        Tnew=[]
        nnew=1
        xnew.append(xmin)
        Tnew.append(T[0])
        while(xnew[nnew-1] < xmax-hmin):</pre>
             for i in range(0,NX-1):
                 if(xnew[nnew-1] >= x[i] and xnew[nnew-1] <= x[i+1] and_{L}
 ⇒xnew[nnew-1]<xmax-hmin):
 _{\text{hll}}=(\text{hloc}[i]*(x[i+1]-\text{xnew}[nnew-1])+\text{hloc}[i+1]*(xnew[nnew-1]-x[i]))/
 \hookrightarrow (x[i+1]-x[i])
                     hll=min(max(hmin,hll),hmax)
                      nnew+=1
                       print(nnew,hll,min(xmax,xnew[nnew-2]+hll))
#
                      xnew.append(min(xmax,xnew[nnew-2]+hll))
#solution interpolation for initialization (attention initial solution on first \Box
 →mesh in the row)
                      un=(T[i]*(x[i+1]-xnew[nnew-1])+T[i+1]*(xnew[nnew-1]-x[i]))/
 \hookrightarrow (x[i+1]-x[i])
                      Tnew.append(un)
        NXO=NX
        NX=nnew
        x = np.zeros((NX))
        x[0:NX] = xnew[0:NX]
        T = np.zeros((NX))
        T[0:NX] = Tnew[0:NX]
```

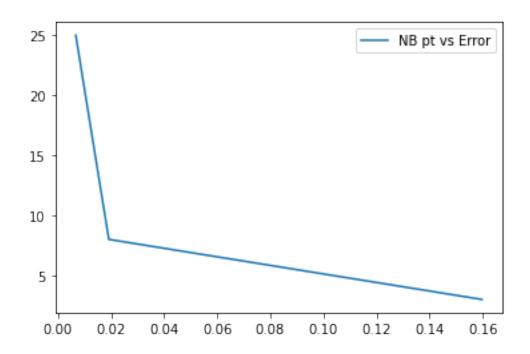
```
T[NX-1]=0
  rest = []
  F = np.zeros((NX))
  RHS = np.zeros((NX))
  hloc = np.ones((NX))*hmax*0.5
  metric = np.ones((NX))
  Tex = np.zeros((NX))
  for j in range (1,NX-1):
      Tex[j] = 2*np.exp(-100*(x[j]-(xmax+xmin)*0.25)**2)+np.
\Rightarrowexp(-200*(x[j]-(xmax+xmin)*0.65)**2)
  dt=1.e30
  for j in range (1,NX-1):
      Tx=(Tex[j+1]-Tex[j-1])/(x[j+1]-x[j-1])
      Txip1=(Tex[j+1]-Tex[j])/(x[j+1]-x[j])
      Txim1=(Tex[j]-Tex[j-1])/(x[j]-x[j-1])
      Txx=(Txip1-Txim1)/(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))
      F[j]=V*Tx-K*Txx+lamda*Tex[j]
      dt=min(dt,0.5*(x[j+1]-x[j-1])**2/(V*np.abs(x[j+1]-x[j-1])+4*K+np.
\Rightarrowabs(F[j])*(x[j+1]-x[j-1])**2))
  print('NX=',NX,'Dt=',dt)
  if(iplot==1):
      plt.figure(1)
  #time step loop
  n=0
  res=1
  res0=1
  t=0
  while(n<NT and res/res0>eps and t<Time):</pre>
      n+=1
      t+=dt
  #discretization of the advection/diffusion/reaction/source equation
      for j in range (1, NX-1):
           visnum=0.5*(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*np.abs(V)
           xnu=K+visnum
          Tx=(T[j+1]-T[j-1])/(x[j+1]-x[j-1])
          Txip1=(T[j+1]-T[j])/(x[j+1]-x[j])
          Txim1=(T[j]-T[j-1])/(x[j]-x[j-1])
          Txx=(Txip1-Txim1)/(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))
          RHS[j] = dt*(-V*Tx+xnu*Txx-lamda*T[j]+F[j])
           metric[j]=min(1./hmin**2,max(1./hmax**2,abs(Txx)/err))
```

```
res+=abs(RHS[j])
      metric[0]=metric[1]
      metric[NX-1]=metric[NX-2] #ux a droite = 0
      \#metric[NX-1]=2*metric[NX-2]-metric[NX-3] \#uxx a droite=0
      for j in range (0, NX-1):
          metric[j]=0.5*(metric[j]+metric[j+1])
      metric[NX-1] = metric[NX-2]
      hloc[0:NX]=np.sqrt(1./metric[0:NX])
      for j in range (1, NX-1):
          T[j] += RHS[j]
          RHS[j]=0
      T[NX-1]=T[NX-2] #1.2*T[NX-2]-0.2*T[NX-3]
      if (n == 1):
          res0=res
      rest.append(res)
  #Plot every ifre time steps
      if (n\%ifre == 0 \text{ or } (res/res0) < eps):
          print('iter=',n,'residual=',res)
          if(iplot==1):
              plotlabel = "t = %1.2f" %(n * dt)
              plt.plot(x[0:NX],T[0:NX], label=plotlabel,linestyle='--',__
→marker='o', color='b')
  print('iter=',n,'time=',t,'residual=',res)
  if(iplot==1):
      plt.plot(x[0:NX],T[0:NX],marker='o', color='b')
      plt.plot(x[0:NX],Tex[0:NX],color='r')
      plt.xlabel(u'$x$', fontsize=26)
      plt.ylabel(u'$T$', fontsize=26, rotation=0)
      plt.title(u'ADRS 1D')
      plt.legend()
      plt.figure(2)
      plt.plot(np.log10(rest/rest[0]))
   errL2=np.sqrt(np.dot(T-Tex,T-Tex))
  errH1h=0
  errL2h=0
```

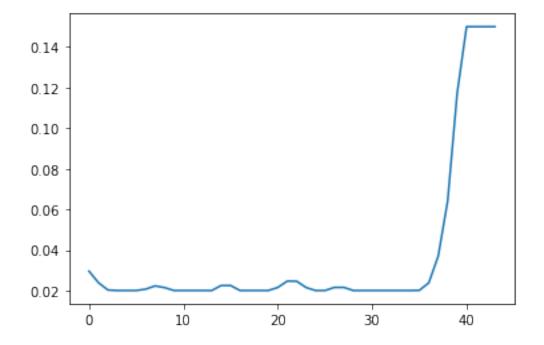
```
for j in range (1, NX-1):
        Texx=(Tex[j+1]-Tex[j-1])/(x[j+1]-x[j-1])
        Tx=(T[j+1]-T[j-1])/(x[j+1]-x[j-1])
        errL2h += (0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*(T[j]-Tex[j])**2
        errH1h+=(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*(Tx-Texx)**2
    errorL2[itera]=errL2h
    errorH1[itera]=errL2h+errH1h
    print('norm error L2, H1=',errL2h,errH1h)
    print('----')
if(iplot==-1):
    plt.figure(3)
    plt.plot(itertab,np.log10(errorL2))
    plt.plot(itertab,np.log10(errorH1))
plt.show()
plt.plot(errorL2[1:itera],1/itertab[1:itera],label="NB pt vs Error")
plt.legend()
plt.show()
plt.plot(hloc)
NX= 8 Dt= 0.05172084342023556
iter= 33 residual= 0.0006021986455240495
iter= 33 time= 1.7067878328677741 residual= 0.0006021986455240495
norm error L2, H1= 0.15967409082372994 2.1050625367114155
-----
NX= 25 Dt= 0.01180541602068137
iter= 112 residual= 0.0016841739743769213
iter= 112 time= 1.322206594316313 residual= 0.0016841739743769213
norm error L2, H1= 0.019131897048811082 4.588746719770155
_____
NX= 43 Dt= 0.007634889695081362
iter= 163 residual= 0.0019908528342445124
iter= 163 time= 1.2444870202982634 residual= 0.0019908528342445124
norm error L2, H1= 0.006671865858585013 2.1428120091486074
NX= 44 Dt= 0.007898086678036072
iter= 158 residual= 0.0021023976231632994
iter= 158 time= 1.2478976951296974 residual= 0.0021023976231632994
norm error L2, H1= 0.006170518730098248 2.0272117740400306
```







[13]: [<matplotlib.lines.Line2D at 0x774764a72d10>]



3.0.1 Premier Workflow

```
[21]: import math
      import numpy as np
      import matplotlib.pyplot as plt
      def adrs_fct(n, x):
      \#u, t = -V u, x + k u, xx - lamda u + f
         u=np.ones(n)
          return u
      def metric_fct(n, u):
      #calcul metric hloc
          hloc=np.ones(n)
          return hloc
      def mesh fct(n, hloc):
      #calcul metric hloc
          x=np.ones(n)
          return x
      iplot=0
      # PHYSICAL PARAMETERS
      K = 0.01 #Diffusion coefficient
      xmin = 0.0
      xmax = 1.0
      Time = 10. #Integration time
      V=1.
      lamda=1
      # ---- (ajout) : background mesh + critère de Cauchy ----
      xB = np.linspace(xmin, xmax, 4000)
      tol_cauchy = 1e-3
      YB_prev = None
      # ---- (ajout) : figure pour tracé multi-étapes -----
      show_progress = True
      plt.figure("Evolution T(x) par adaptation - WF1")
      # mesh adaptation param
      niter_refinement=30
      hmin=0.02
      hmax=0.15
      err=0.03
```

```
# NUMERICAL PARAMETERS
NX = 3
NT = 10000
ifre=1000000
eps=0.001
errorL2=np.zeros((niter_refinement))
errorH1=np.zeros((niter refinement))
itertab=np.zeros((niter_refinement))
hloc = np.ones((NX))*hmax
itera=0
NXO=0
while( np.abs(NXO-NX) > 2 and itera<niter_refinement-1):</pre>
    itera+=1
    itertab[itera]=1./NX
    iplot=itera-2
    x = np.linspace(xmin,xmax,NX)
    T = np.zeros((NX))
    # mesh adaptation using local metric
    if(itera>0):
        xnew=[]
        Tnew=[]
        nnew=1
        xnew.append(xmin)
        Tnew.append(T[0])
        while(xnew[nnew-1] < xmax-hmin):</pre>
             for i in range(0,NX-1):
                  if(xnew[nnew-1] >= x[i] and xnew[nnew-1] <= x[i+1] and_{\sqcup}
 →xnew[nnew-1]<xmax-hmin):</pre>
 _{\text{hll}}=(\text{hloc}[i]*(x[i+1]-\text{xnew}[nnew-1])+\text{hloc}[i+1]*(xnew[nnew-1]-x[i]))/
 \hookrightarrow (x[i+1]-x[i])
                      hll=min(max(hmin,hll),hmax)
                      nnew+=1
                      xnew.append(min(xmax,xnew[nnew-2]+hll))
                      # interpolation de T (initialisation)
                      un=(T[i]*(x[i+1]-xnew[nnew-1])+T[i+1]*(xnew[nnew-1]-x[i]))/
 \hookrightarrow (x[i+1]-x[i])
                      Tnew.append(un)
        NXO=NX
```

```
NX=nnew
      x = np.zeros((NX))
      x[0:NX] = xnew[0:NX]
      T = np.zeros((NX))
      T[0:NX] = Tnew[0:NX]
  rest = []
  F = np.zeros((NX))
  RHS = np.zeros((NX))
  hloc = np.ones((NX))*hmax*0.5
  metric = np.ones((NX))
  Tex = np.zeros((NX))
  for j in range (1,NX-1):
      Tex[j] = 2*np.exp(-100*(x[j]-(xmax+xmin)*0.25)**2)+np.
\Rightarrowexp(-200*(x[j]-(xmax+xmin)*0.65)**2)
  dt=1.e30
  for j in range (1,NX-1):
      Tx=(Tex[j+1]-Tex[j-1])/(x[j+1]-x[j-1])
      Txip1=(Tex[j+1]-Tex[j])/(x[j+1]-x[j])
      Txim1 = (Tex[j] - Tex[j-1])/(x[j] - x[j-1])
      Txx=(Txip1-Txim1)/(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))
      F[j]=V*Tx-K*Txx+lamda*Tex[j]
      dt=min(dt,0.5*(x[j+1]-x[j-1])**2/(V*np.abs(x[j+1]-x[j-1])+4*K+np.
\Rightarrowabs(F[j])*(x[j+1]-x[j-1])**2))
  print('NX=',NX,'Dt=',dt)
  if(iplot==1):
      plt.figure(1)
  # time step loop
  n=0
  res=1
  res0=1
  t=0
  while(n<NT and res/res0>eps and t<Time):</pre>
      n+=1
      t+=dt
      res=0
      for j in range (1, NX-1):
           visnum=0.5*(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*np.abs(V)
           xnu=K+visnum
           Tx=(T[j+1]-T[j-1])/(x[j+1]-x[j-1])
           Txip1=(T[j+1]-T[j])/(x[j+1]-x[j])
           Txim1=(T[j]-T[j-1])/(x[j]-x[j-1])
```

```
Txx = (Txip1 - Txim1) / (0.5*(x[j+1] + x[j]) - 0.5*(x[j] + x[j-1]))
           RHS[j] = dt*(-V*Tx+xnu*Txx-lamda*T[j]+F[j])
          metric[j]=min(1./hmin**2,max(1./hmax**2,abs(Txx)/err))
           res+=abs(RHS[j])
      metric[0] = metric[1]
      metric[NX-1]=metric[NX-2]
      for j in range (0, NX-1):
          metric[j]=0.5*(metric[j]+metric[j+1])
      metric[NX-1]=metric[NX-2]
      hloc[0:NX]=np.sqrt(1./metric[0:NX])
      for j in range (1, NX-1):
          T[j] += RHS[j]
          RHS[j]=0
      T[NX-1]=T[NX-2]
      if (n == 1):
          res0=res
      rest.append(res)
      if (n\%ifre == 0 \text{ or } (res/res0) < eps):
           print('iter=',n,'residual=',res)
           if(iplot==1):
               plotlabel = "t = %1.2f" %(n * dt)
               plt.plot(x[0:NX],T[0:NX], label=plotlabel,linestyle='--',_u
→marker='o', color='b')
  print('iter=',n,'time=',t,'residual=',res)
  # ---- (ajout) : tracer la solution à chaque adaptation ----
  if show_progress:
      plt.plot(x, T, '-', alpha=0.5)
  # ---- projection background + Cauchy (WF1) ----
  YB = np.interp(xB, x, T)
  if YB_prev is not None:
      Ik = np.sqrt(np.trapz((YB - YB_prev)**2, xB))
      print(f'Cauchy Ik = {Ik:.3e}')
      if Ik < tol_cauchy:</pre>
           print(' Critère de Cauchy atteint : arrêt de l'adaptation.')
           break
```

```
YB_prev = YB.copy()
    if(iplot==1):
       plt.plot(x[0:NX],T[0:NX],marker='o', color='b')
       plt.plot(x[0:NX],Tex[0:NX],color='r')
       plt.xlabel(u'$x$', fontsize=26)
       plt.ylabel(u'$T$', fontsize=26, rotation=0)
       plt.title(u'ADRS 1D')
       plt.legend()
       plt.figure(2)
       plt.plot(np.log10(rest/rest[0]))
   errH1h=0
   errL2h=0
   for j in range (1, NX-1):
       Texx=(Tex[j+1]-Tex[j-1])/(x[j+1]-x[j-1])
       Tx=(T[j+1]-T[j-1])/(x[j+1]-x[j-1])
       errL2h += (0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*(T[j]-Tex[j])**2
        errH1h+=(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*(Tx-Texx)**2
   errorL2[itera]=errL2h
   errorH1[itera]=errL2h+errH1h
   print('norm error L2, H1=',errL2h,errH1h)
   print('----')
# --- (ajout) : finaliser la figure d'évolution ----
if show_progress:
   plt.xlabel("x"); plt.ylabel("T")
   plt.title("ADRS: évolution par adaptation (workflow 1)")
   plt.grid(True, alpha=0.3)
if(iplot==-1):
   plt.figure(3)
   plt.plot(itertab,np.log10(errorL2))
   plt.plot(itertab,np.log10(errorH1))
plt.show()
plt.plot(errorL2[1:itera],1/itertab[1:itera],label="NB pt vs Error")
plt.legend()
plt.show()
plt.plot(hloc)
```

NX= 8 Dt= 0.05172084342023556

iter= 33 residual= 0.0006021986455240495
iter= 33 time= 1.7067878328677741 residual= 0.0006021986455240495
norm error L2, H1= 0.15967409082372994 2.1050625367114155

NX= 25 Dt= 0.01180541602068137

iter= 112 residual= 0.0016841739743769213

iter= 112 time= 1.322206594316313 residual= 0.0016841739743769213

Cauchy Ik = 4.184e-01

norm error L2, H1= 0.019131897048811082 4.588746719770155

NX= 43 Dt= 0.007634889695081362

iter= 163 residual= 0.0019908528342445124

iter= 163 time= 1.2444870202982634 residual= 0.0019908528342445124

Cauchy Ik = 7.020e-02

norm error L2, H1= 0.006671865858585013 2.1428120091486074

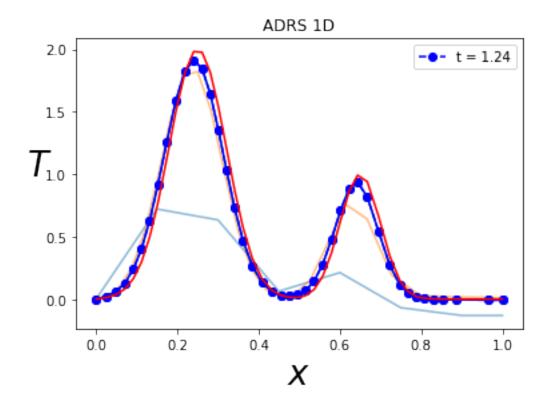
NX= 44 Dt= 0.007898086678036072

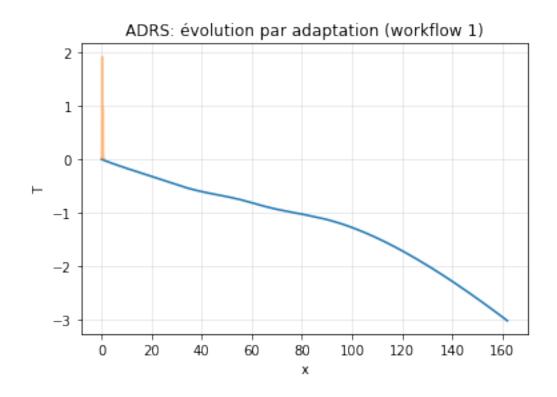
iter= 158 residual= 0.0021023976231632994

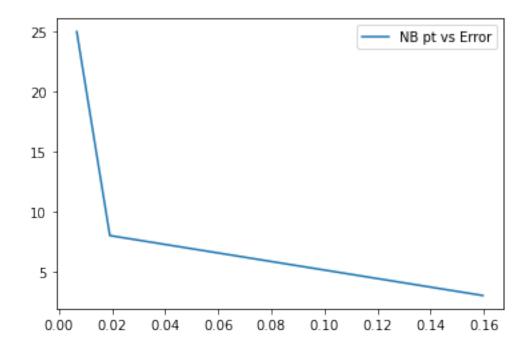
iter= 158 time= 1.2478976951296974 residual= 0.0021023976231632994

Cauchy Ik = 1.289e-02

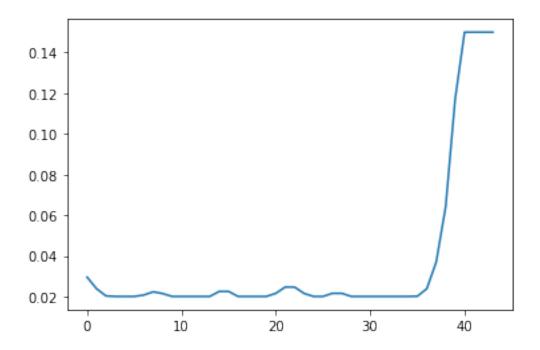
norm error L2, H1= 0.006170518730098248 2.0272117740400306







[21]: [<matplotlib.lines.Line2D at 0x774765069d50>]



3.0.2 Second Workflow

```
[25]: import math
      import numpy as np
      import matplotlib.pyplot as plt
      def adrs_fct(n, x):
      \#u, t = -V u, x + k u, xx - lamda u + f
          u=np.ones(n)
          return u
      def metric_fct(n, u):
      #calcul metric hloc
          hloc=np.ones(n)
          return hloc
      def mesh_fct(n, hloc):
      #calcul metric hloc
          x=np.ones(n)
          return x
      iplot=0
      # PHYSICAL PARAMETERS
      K = 0.01
                   \#Diffusion\ coefficient
      xmin = 0.0
```

```
xmax = 1.0
Time = 10. #Integration time
V=1.
lamda=1
# >>> MODIF (workflow 2): background mesh + tolérance Cauchy
xB = np.linspace(xmin, xmax, 4000)
tol_cauchy = 1e-3
YB_prev = None
# >>> MODIF (tracé multi-étapes, minimal)
show_progress = True
plt.figure("Evolution T(x) par adaptation - WF2")
#mesh adaptation param
niter_refinement=30 #niter different calculations
hmin=0.02
hmax=0.15
err=0.03
# NUMERICAL PARAMETERS
NX = 3 #Number of grid points : initialization
NT = 10000 #Number of time steps max
ifre=1000000 #plot every ifre time iterations
eps=0.001 #relative convergence ratio
errorL2=np.zeros((niter_refinement))
errorH1=np.zeros((niter_refinement))
itertab=np.zeros((niter_refinement))
\# >>> MODIF: initialisation unique (x,T) AVANT la boucle
x = np.linspace(xmin, xmax, NX)
T = np.zeros((NX))
# -----
hloc = np.ones((NX))*hmax
itera=0
NXO=0
while( np.abs(NXO-NX) > 2 and itera<niter_refinement-1):</pre>
   itera+=1
   itertab[itera]=1./NX
   iplot=itera-2
```

```
# >>> MODIF: NE PAS re-créer x, T ici (on garde ceux de l'itération_{\sqcup}
⇔précédente)
  \# x = np.linspace(xmin, xmax, NX)
   \# T = np.zeros((NX))
   # >>> MODIF: conserver l'état avant remeshing pour interpolation_
\hookrightarrow Y_{k-1}-Mesh_{k}
  x_old = x.copy()
  T_old = T.copy()
   #mesh adaptation using local metric
  if(itera>0):
       xnew=[]
       Tnew=[]
       nnew=1
       xnew.append(xmin)
       # >>> MODIF: partir de la vraie solution précédente
       Tnew.append(T_old[0])
       # -----
       while(xnew[nnew-1] < xmax-hmin):</pre>
           for i in range(0,NX-1):
                if(xnew[nnew-1] >= x[i] and xnew[nnew-1] <= x[i+1] and_{\sqcup}
→xnew[nnew-1]<xmax-hmin):</pre>
_{\circ}hll=(hloc[i]*(x[i+1]-xnew[nnew-1])+hloc[i+1]*(xnew[nnew-1]-x[i]))/
\hookrightarrow (x[i+1]-x[i])
                    hll=min(max(hmin,hll),hmax)
                    nnew+=1
                    xnew.append(min(xmax,xnew[nnew-2]+hll))
                    # >>> MODIF: interpolation de Y_{k-1} (T_{old}) sur le
⇔nouveau noeud
                    un=(T_old[i]*(x[i+1]-xnew[nnew-1]) +_{\square}
\neg T_old[i+1]*(xnew[nnew-1]-x[i]))/(x[i+1]-x[i])
                    Tnew.append(un)
       NXO=NX
       NX=nnew
       # >>> MODIF: ne PAS réinitialiser T à zéro; on prend l'interpolation
       x = np.array(xnew[:NX])
       T = np.array(Tnew[:NX])
```

```
# T[NX-1]=0 # (laisse commenté)
  rest = []
  F = np.zeros((NX))
  RHS = np.zeros((NX))
  hloc = np.ones((NX))*hmax*0.5
  metric = np.ones((NX))
  Tex = np.zeros((NX))
  for j in range (1,NX-1):
      Tex[j] = 2*np.exp(-100*(x[j]-(xmax+xmin)*0.25)**2)+np.
\Rightarrowexp(-200*(x[j]-(xmax+xmin)*0.65)**2)
  dt=1.e30
  for j in range (1,NX-1):
      Tx=(Tex[j+1]-Tex[j-1])/(x[j+1]-x[j-1])
      Txip1=(Tex[j+1]-Tex[j])/(x[j+1]-x[j])
      Txim1 = (Tex[j] - Tex[j-1])/(x[j] - x[j-1])
      Txx=(Txip1-Txim1)/(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))
      F[j]=V*Tx-K*Txx+lamda*Tex[j]
      dt=min(dt,0.5*(x[j+1]-x[j-1])**2/(V*np.abs(x[j+1]-x[j-1])+4*K+np.
\Rightarrowabs(F[j])*(x[j+1]-x[j-1])**2))
  print('NX=',NX,'Dt=',dt)
  if(iplot==1):
      plt.figure(1)
  #time step loop (SOLVE sur Mesh_k en repartant de l'interpolation Y_{k-1})
  n=0
  res=1
  res0=1
  t.=()
  while(n<NT and res/res0>eps and t<Time):</pre>
      n+=1
      t+=dt
      #discretization of the advection/diffusion/reaction/source equation
      for j in range (1, NX-1):
          visnum=0.5*(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*np.abs(V)
           xnu=K+visnum
          Tx=(T[j+1]-T[j-1])/(x[j+1]-x[j-1])
          Txip1=(T[j+1]-T[j])/(x[j+1]-x[j])
          Txim1=(T[j]-T[j-1])/(x[j]-x[j-1])
          Txx=(Txip1-Txim1)/(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))
          RHS[j] = dt*(-V*Tx+xnu*Txx-lamda*T[j]+F[j])
           metric[j]=min(1./hmin**2,max(1./hmax**2,abs(Txx)/err))
```

```
res+=abs(RHS[j])
      metric[0] = metric[1]
      metric[NX-1]=metric[NX-2] #ux a droite = 0
      for j in range (0, NX-1):
          metric[j]=0.5*(metric[j]+metric[j+1])
      metric[NX-1] = metric[NX-2]
      hloc[0:NX]=np.sqrt(1./metric[0:NX])
      for j in range (1, NX-1):
           T[j] += RHS[j]
          RHS[j]=0
      T[NX-1]=T[NX-2] #1.2*T[NX-2]-0.2*T[NX-3]
      if (n == 1):
          res0=res
      rest.append(res)
      if (n\%ifre == 0 \text{ or } (res/res0) < eps):
          print('iter=',n,'residual=',res)
           if(iplot==1):
               plotlabel = "t = %1.2f" %(n * dt)
               plt.plot(x[0:NX],T[0:NX], label=plotlabel,linestyle='--',__

marker='o', color='b')
  print('iter=',n,'time=',t,'residual=',res)
  # >>> MODIF (tracé à chaque adaptation, minimal)
  if show_progress:
      plt.plot(x, T, '-', alpha=0.55)
  # >>> MODIF (workflow 2): projection sur background + test de Cauchy
  YB = np.interp(xB, x, T)
  if YB_prev is not None:
      Ik = np.sqrt(np.trapz((YB - YB_prev)**2, xB))
      print(f'Cauchy Ik = {Ik:.3e}')
      if Ik < tol_cauchy:</pre>
           print(' Critère de Cauchy atteint : arrêt de l'adaptation.')
          break
  YB_prev = YB.copy()
```

```
if(iplot==1):
        plt.plot(x[0:NX],T[0:NX],marker='o', color='b')
        plt.plot(x[0:NX],Tex[0:NX],color='r')
        plt.xlabel(u'$x$', fontsize=26)
        plt.ylabel(u'$T$', fontsize=26, rotation=0)
        plt.title(u'ADRS 1D')
        plt.legend()
        plt.figure(2)
        plt.plot(np.log10(rest/rest[0]))
    errH1h=0
    errL2h=0
    for j in range (1, NX-1):
        Texx=(Tex[j+1]-Tex[j-1])/(x[j+1]-x[j-1])
        Tx=(T[j+1]-T[j-1])/(x[j+1]-x[j-1])
        errL2h+=(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*(T[j]-Tex[j])**2
        errH1h+=(0.5*(x[j+1]+x[j])-0.5*(x[j]+x[j-1]))*(Tx-Texx)**2
    errorL2[itera]=errL2h
    errorH1[itera]=errL2h+errH1h
    print('norm error L2, H1=',errL2h,errH1h)
    print('----')
# >>> MODIF (finalisation de la figure)
if show_progress:
    plt.xlabel("x"); plt.ylabel("T")
    plt.title("ADRS: évolution par adaptation (workflow 2)")
    plt.grid(True, alpha=0.3)
    plt.show()
NX= 8 Dt= 0.05172084342023556
iter= 33 residual= 0.0006021986455240495
iter= 33 time= 1.7067878328677741 residual= 0.0006021986455240495
norm error L2, H1= 0.15967409082372994 2.1050625367114155
-----
NX= 25 Dt= 0.01178864148943053
iter= 97 residual= 0.0014715386835857688
iter= 97 time= 1.1434982244747607 residual= 0.0014715386835857688
Cauchy Ik = 4.211e-01
norm error L2, H1= 0.018085733097822607 4.522077151124964
NX= 40 Dt= 0.007430241465423633
iter= 108 residual= 0.0010563628614907168
iter= 108 time= 0.802466078265751 residual= 0.0010563628614907168
Cauchy Ik = 6.932e-02
norm error L2, H1= 0.006385254501066739 2.084604123692186
```

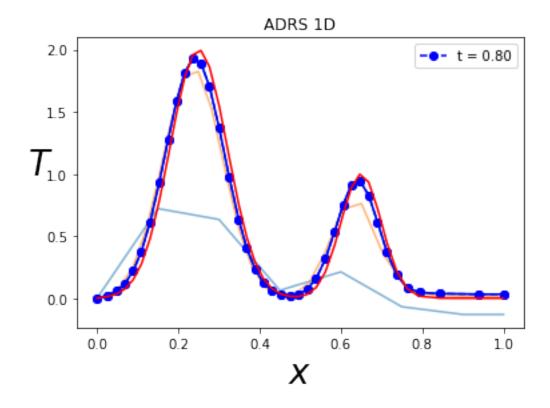
NX= 40 Dt= 0.0077083056771561605

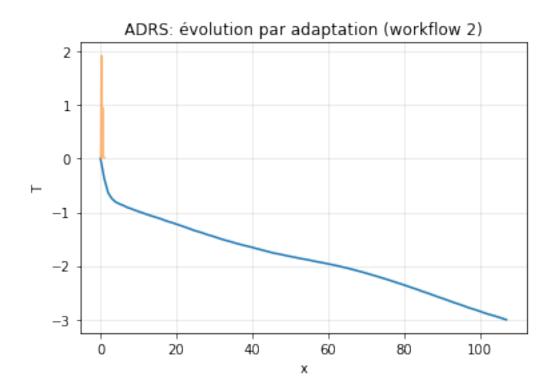
iter= 102 residual= 0.0002725602090546944

iter= 102 time= 0.7862471790699288 residual= 0.0002725602090546944

Cauchy Ik = 9.164e-03

norm error L2, H1= 0.006073746151591898 1.9589484836332234





[]: