Cosmology: Project 1.

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Point 1:

Find solutions for the scale factor a(t) using the cosmological parameters of the baseline cosmology, use it to express the Hubble factor H as function of t.

General form:

$$rac{H^2}{H_o^2} = rac{\dot{a}^2}{a^2 H_0^2} = rac{\Omega_{r,0}}{a^4} + rac{\Omega_{m,0}}{a^3} + \Omega_{\Lambda,0} + rac{1-\Omega_0}{a^2}$$

The benchmark model has $\Omega_0=1$, then we have:

$$\implies \dot{a} = H_0 \sqrt{rac{\Omega_{r,0}}{a^2} + rac{\Omega_{m,0}}{a} + \Omega_{\Lambda,0} a^2}$$

```
import the necessary packages
import numpy as np # Package for scientific computing with Python
import matplotlib.pyplot as plt # Package for plotting
import pandas as pd # Package for data analysis
import seaborn as sns # Package for plotting styles
import matplotlib as mlp # Package for plotting

sns.set_theme()

# Improve image resolution
mlp.rcParams['figure.dpi'] = 100
```

```
In [2]: # Important constants
    c = 299792.458 # Units: km/s
    G = 4.302e-9 # Units: km^2 Mpc^-3 Msun^-1 s^-2

# Cosmological parameters from Planck 2015 results
    H_0 = 67.36 # Units: km/s/Mpc
    Omega_m = 0.3111 # Matter density
    Omega_l = 0.6889 # Dark energy density
    Omega_r = 4.165e-5 # Radiation density
```

```
INPUT:
    a: expansion factor
    radiation: if True, include radiation
    matter: if True, include matter
    dark: if True, include dark energy
OUTPUT:
    a_dot: derivative of the expansion factor
# Contribution from different types of matter
a m = omega m/a
a r = omega r/(a**2)
a_1 = omega_1*(a**2)
a_dot = a_m + a_l + a_r
# Check for equality
if equality:
    if np.abs(a_m - a_1) < 0.0001 and a_m != 0 and a_1 != 0:</pre>
        print('Matter and dark matter equality at a = ', a)
    if np.abs(a m - a r) < 0.0001 and a m != 0 and a r != 0:</pre>
        print('Matter and radiation equality at a = ', a)
    if np.abs(a l - a r) < 0.0001 and a l != 0 and a r != 0:</pre>
        print('Dark matter and radiation equality at a = ', a)
# Not negative values allowed
if a dot <0:</pre>
    print('Negative derivative at a = ', a)
    return 0
# Return the derivative
a_dot = H_0 * np.sqrt(a_dot)
return a dot
```

```
In [4]:
         def runge kutta(f, y 0, t0 = 0, h = -0.001, omega m = Omega m,
                         omega 1 = Omega 1, omega r = Omega r, return densities = False, n max =
             Function to solve a system of differential equations using the Runge-Kutta method (
             INPUT:
                 f: function to solve
                 t0: initial time
                 y_0: initial value
                 h: step
             OUTPUT:
                 t: time
                 y: solution
             # Initialize list to store densities
             matter, radiation, dark = [omega_m], [omega_r], [omega_1]
             # Initialize list to store a derivatives
             a_dot = [get_a_dot(y_0, omega_m = omega_m, omega_l = omega_l, omega_r = omega_r)]
             # Initialize list to store time and a.
             y = [y_0]
             t = [t0]
             # Loop to solve the system of differential equations
             while True:
```

```
# Runge-Kutta coefficients
        k1 = f(y[-1], omega_m = omega_m, omega_l = omega_l, omega_r = omega_r)
        k2 = f(y[-1] + k1*h/2, omega_m = omega_m, omega_l = omega_l, omega_r = omega_r)
        k3 = f(y[-1] + k2*h/2, omega_m = omega_m, omega_l = omega_l, omega_r = omega_r)
        k4 = f(y[-1] + k3*h, omega_m = omega_m, omega_1 = omega_1, omega_r = omega_r)
        # Update time and solution
        y_aux = y[-1] + (k1 + 2*k2 + 2*k3 + k4)*h/6
        t_{aux} = t[-1] + h
        # Maximum number of iterations
        if len(y) > n_max:
            break
        # Only allow positive a and lower than 20
        if (h < 0 and y_aux > 0) or (h > 0 and y_aux < 20):</pre>
                y.append(y_aux)
                t.append(t_aux)
                # If return densities is True, store the densities
                if return_densities:
                    # Store the derivatives
                    y_{dot} = f(y_{aux}, omega_m = omega_m, omega_1 = omega_1, omega_r = omega_n
                    # Compute H 0/H
                    constant = H_0**2/(y_dot/y_aux)**2
                    # Store the densities
                    a_dot.append(y_dot)
                    matter.append(omega_m/(y_aux**3)*constant)
                    radiation.append(omega_r/(y_aux**4)*constant)
                    dark.append(omega_l*constant)
        else:
            break
    if return_densities:
        return t, y, matter, radiation, dark, a_dot
    else:
        return t, y
h = 0.001 # Step
```

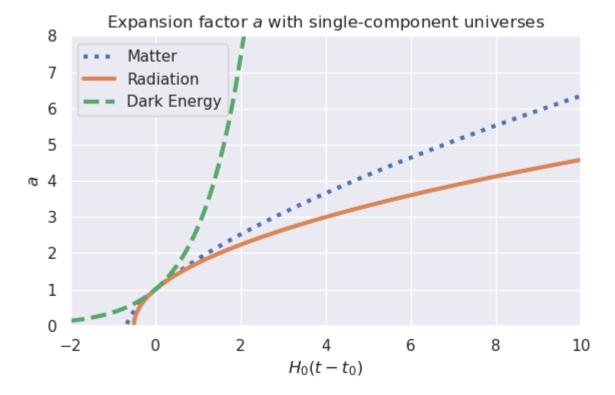
```
In [5]:
         t0 = 1 # Initial time
         # Universe only with matter
         x_m, y_m = runge_kutta(f = get_a_dot, y_0 = 1,t0 = t0, h = h,
                             omega_m = 1.0, omega_l = 0.0, omega_r = 0.0) # Future expansion fac
         x_{aux}, y_{aux} = runge_kutta(f = get_a_dot, y_0 = 1, t0 = t0, h = -0.0001,
                              omega_m = 1.0, omega_l = 0.0, omega_r = 0.0) # Past expansion facto
         # Merge the two lists
         x_m = x_aux[::-1] + x_m
         y_m = y_aux[::-1] + y_m
         # Universe only with radiation
         x_r, y_r = runge_kutta(f = get_a_dot, y_0 = 1, t0 = t0, h = h, omega_m = 0.0, omega_1
         x_{aux}, y_{aux} = runge_kutta(f = get_a_dot, y_0 = 1, t0 = t0, h = -0.0001, omega_m = 0.0)
         x_r = x_{aux}[::-1] + x_r
         y_r = y_aux[::-1] + y_r
         # Universe only with dark matter
```

```
x_1, y_1 = runge_kutta(f = get_a_dot, y_0 = 1, t0 = t0, h = h, omega_m = 0.0, omega_1
x_aux, y_aux = runge_kutta(f = get_a_dot, y_0 = 1, t0 = t0, h = -h, omega_m = 0.0, omeg
x_1 = x_aux[::-1] + x_1
y_1 = y_aux[::-1] + y_1
```

/tmp/ipykernel_543/3429811448.py:15: RuntimeWarning: invalid value encountered in double
_scalars

```
a_r = omega_r/(a^{**2})
```

```
In [6]:
         # Plot of single-component universes
         fig, ax = plt.subplots(figsize = (6, 4))
         ax.plot((np.array(x_m)-t0)*H_0, y_m, label = 'Matter', linestyle = 'dotted', linewidth
         ax.plot((np.array(x_r)-t0)*H_0, y_r, label = 'Radiation', linestyle = 'solid', linewidt')
         ax.plot((np.array(x_1)-t0)*H_0, y_1, label = 'Dark Energy', linestyle = 'dashed', linew
         ax.set_ylim(0, 8)
         ax.set xlim(-2, 10)
         ax.set_xlabel('$H_0(t-t_0)$')
         ax.set ylabel('$a$')
         ax.set_title('Expansion factor $a$ with single-component universes')
         ax.legend()
         plt.tight layout()
         plt.savefig('./figs/single_component_universe.png', dpi = 300)
         plt.show()
         plt.close(fig)
```



```
In [7]: # Simulate benchmark universe model (three components)

t0 = 0
x_b, y_b = runge_kutta(f = get_a_dot, y_0 = 1, t0 = t0, h = 0.0000001, n_max = 1000000)
x_aux, y_aux = runge_kutta(f = get_a_dot, y_0 = 1, t0 = t0, h = -0.0000001, n_max = 100
```

```
x_b = x_{aux}[::-1] + x_b

y_b = y_{aux}[::-1] + y_b
```

```
# Matter-Dark matter equality at 0.7672
idx_ml = np.where(np.abs(np.array(y_b) - 0.76725) < 0.0001)[0][0]
print(x_b[idx_ml], y_b[idx_ml])

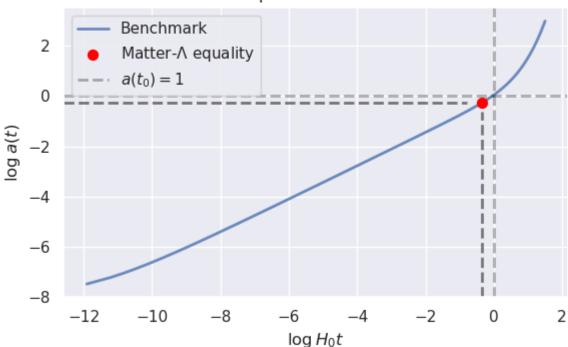
# Radiation-Matter (acording to the book) at 2.9e-4
idx_rm = np.where(np.abs(np.array(y_b) - 2.9e-4) < 1)[0][0]
print(x_b[idx_rm], y_b[idx_rm])</pre>
```

- -0.003658699999997579 0.7671549583080319
- -0.01416469999997014 0.00035275320459171697

```
In [9]:
         # Benchmark model plot (log scale)
         fig, ax = plt.subplots(figsize = (6, 4))
         ax.plot(np.log((np.array(x b)-min(x b))*H 0), np.log(y b), label = 'Benchmark', linesty
         # Matter-Dark matter equality at 0.7672
         x_ml = np.log((np.array(x_b[idx_ml])-min(x_b))*H_0)
         y ml = np.log(y b[idx ml])
         ax.scatter(x_ml, y_ml, color = 'red', label = 'Matter-$\Lambda$ equality', marker = 'o'
         ax.axhline(y = y_ml, color = 'black', linestyle = 'dashed', linewidth = 2, alpha = 0.5,
         ax.axvline(x = x_ml, color = 'black', linestyle = 'dashed', linewidth = 2, alpha = 0.5,
         # Today
         ax.axhline(y = 0, color = 'black', linestyle = 'dashed', alpha = 0.3, linewidth = 2)
         ax.axvline(x = 0, color = 'black', linestyle = 'dashed', alpha = 0.3, linewidth = 2, la
         ax.set xlabel('$\log{H Ot}$')
         ax.set ylabel('$\log{a(t)}$')
         ax.set_title('Expansion factor $a$')
         ax.legend()
         plt.tight layout()
         plt.savefig('./figs/benchmark universe log.png', dpi = 300)
         plt.show()
         plt.close(fig)
```

/tmp/ipykernel_543/3747719866.py:4: RuntimeWarning: divide by zero encountered in log ax.plot(np.log((np.array(x_b)-min(x_b))*H_0), np.log(y_b), label = 'Benchmark', linest yle = 'solid', linewidth = 2, alpha = 0.8)

Expansion factor a

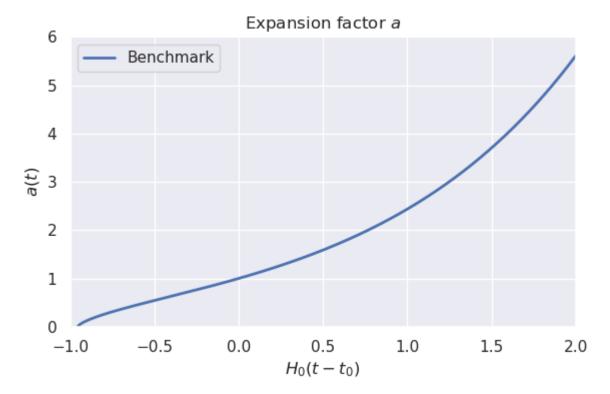


```
In [10]:
# Benchmark model plot (linear scale)
fig, ax = plt.subplots(figsize = (6, 4))

ax.plot((np.array(x_b)-t0)*H_0, y_b, label = 'Benchmark', linestyle = 'solid', linewidt

ax.set_xlabel('$H_0(t-t_0)$')
ax.set_ylabel('$a(t)$')
ax.set_ylim(0, 6)
ax.set_ylim(0, 6)
ax.set_xlim(-1, 2)
ax.set_title('Expansion factor $a$')
ax.legend()

plt.tight_layout()
plt.savefig('./figs/benchmark_universe.png', dpi = 300)
plt.show()
plt.close(fig)
```



Point 2:

Make plots of how the energy density of the different matter-energy species evolves as function of a, z and t.

We know how to calculate the energy density of the different matter-energy species. We also know, from the previous point, how the scale factor changes as a function of time. We will use the benchmark model to calculate the energy density of the different matter-energy species as a function of time.

The relationship between the energy density of radiation with respect to the scale factor a is given by:

$$egin{align} arepsilon_r &= rac{arepsilon_{r,0}}{a^4} \quad ; \quad \Omega_r &= rac{arepsilon_r}{arepsilon_c} \quad ; \quad arepsilon_c &= rac{3c^2}{8\pi G} (rac{\dot{a}}{a})^2 \ &\Longrightarrow \ \Omega_r &= rac{\Omega_{r,0} arepsilon_{c,0}}{a^4 arepsilon_c} \implies \Omega_r &= rac{\Omega_{r,0}}{a^4} rac{H_0^2}{H^2} \ \end{split}$$

Analogously, the energy density of matter and dark energy is given by:

$$\Omega_m = rac{\Omega_{m,0}}{a^3} rac{H_0^2}{H^2} \quad ; \quad \Omega_\Lambda = \Omega_{\Lambda,0} rac{H_0^2}{H^2}$$

```
# Benchmark model with densities

t0 = 0

x_b, y_b, matter, radiation, dark, a_dot = runge_kutta(f = get_a_dot, y_0 = 1,

t0 = t0, h = 0.0000001, return_densities=True, n_max = 1000000)

x aux, y aux, matter aux, radiation aux, dark aux, a dot aux = runge kutta(f = get a do
```

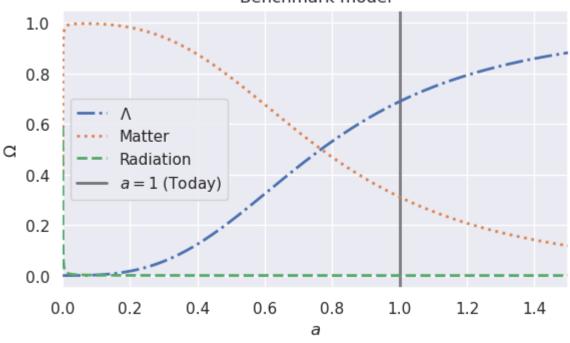
```
h = -0.00000001, return_densities=True, n_max = 10000000)

x_b = x_aux[::-1] + x_b
y_b = y_aux[::-1] + y_b
matter = matter_aux[::-1] + matter
radiation = radiation_aux[::-1] + radiation
dark = dark_aux[::-1] + dark
a_dot = a_dot_aux[::-1] + a_dot
```

```
a matter density radiation density dark density
Out[12]:
                                                                                     a_dot
           0 -0.954140 0.000093
                                       0.409850
                                                        0.590150 7.294826e-13 6086.261121
           1 -0.954139 0.000144
                                       0.517432
                                                        0.482568 3.389521e-12 4359.339198
          2 -0.954138 0.000183
                                       0.577814
                                                        0.422186 7.871165e-12 3651.396884
          3 -0.954138 0.000218
                                       0.619015
                                                        0.380985 1.410864e-11 3237.775637
           4 -0.954137 0.000248
                                       0.649805
                                                        0.350195 2.205987e-11 2957.104080
```

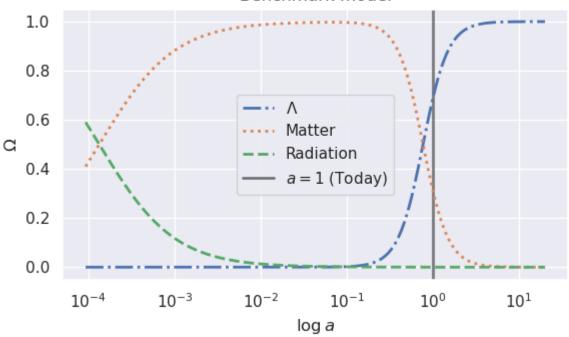
```
In [13]:
          # Densities evolution as a function of the scale factor
          fig, ax = plt.subplots(figsize = (6, 4))
          ax.plot(df['a'], df['dark density'], label = '$\Lambda$', linestyle = 'dashdot', linewi
          ax.plot(df['a'], df['matter density'], label = 'Matter', linestyle = 'dotted', linewidt
          ax.plot(df['a'], df['radiation density'], label = 'Radiation', linestyle = 'dashed', li
          ax.axvline(x = 1, color = 'black', linestyle = 'solid', linewidth = 2, alpha = 0.5, lab
          ax.set xlabel('$a$')
          ax.set ylabel('$\Omega$')
          ax.set title('Evolution of densities as a function of time.\nBenchmark model')
          ax.set xlim(0, 1.5)
          plt.legend(loc = 'best')
          plt.tight layout()
          plt.savefig('./figs/densities evolution a.png', dpi = 300)
          plt.show()
          plt.close(fig)
```

Evolution of densities as a function of time. Benchmark model



```
In [14]:
          # Densities evolution as a function of the scale factor (log scale)
          fig, ax = plt.subplots(figsize = (6, 4))
          ax.plot(df['a'], df['dark density'], label = '$\Lambda$', linestyle = 'dashdot', linewi
          ax.plot(df['a'], df['matter density'], label = 'Matter', linestyle = 'dotted', linewidt
          ax.plot(df['a'], df['radiation density'], label = 'Radiation', linestyle = 'dashed', li
          ax.axvline(x = 1, color = 'black', linestyle = 'solid', linewidth = 2, alpha = 0.5, lab
          ax.set_xlabel('$\log{a}$')
          ax.set ylabel('$\Omega$')
          ax.set_title('Evolution of densities as a function of time.\nBenchmark model')
          # X-scale to log scale
          plt.xscale('log')
          plt.legend(loc = 'center')
          plt.tight layout()
          plt.savefig('./figs/densities_evolution_loga.png', dpi = 300)
          plt.show()
          plt.close(fig)
```

Evolution of densities as a function of time. Benchmark model



The redshift is related with the scale factor as:

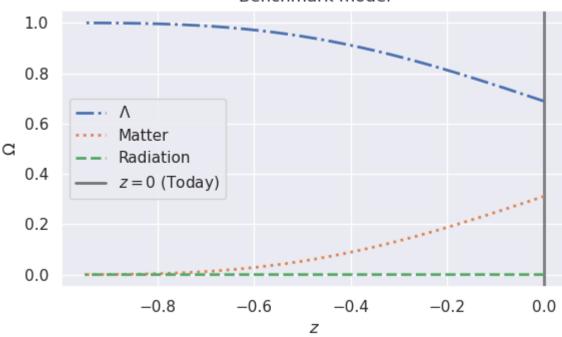
$$1+z=\frac{1}{a(t)}$$

With $a(t_{obs}) = 1$.

```
In [15]: # Get redshift for every expansion factor
df['z'] = 1/df['a'] - 1
```

```
In [16]:
          # Denisties evolution as a function of redshift (z<0 - Future)
          fig, ax = plt.subplots(figsize = (6, 4))
          df_future = df[df['z'] < 0]</pre>
          df past = df[df['z'] > 0.01]
          ax.plot(df_future['z'], df_future['dark density'], label = '$\Lambda$', linestyle = 'da
          ax.plot(df_future['z'], df_future['matter density'], label = 'Matter', linestyle = 'dot
          ax.plot(df future['z'], df future['radiation density'], label = 'Radiation', linestyle
          ax.axvline(x = 0, color = 'black', linestyle = 'solid', linewidth = 2, alpha = 0.5, lab
          ax.set_xlabel('$z$')
          ax.set ylabel('$\Omega$')
          ax.set title('Evolution of densities as a function of $z<0$.\nBenchmark model')</pre>
          plt.legend(loc = 'best')
          plt.tight layout()
          plt.savefig('./figs/densities evolution z future.png', dpi = 300)
          plt.show()
          plt.close(fig)
```

Evolution of densities as a function of z < 0. Benchmark model



```
In [17]:
# Denisties evolution as a function of redshift (z>0 - past)
fig, ax = plt.subplots(figsize = (6, 4))

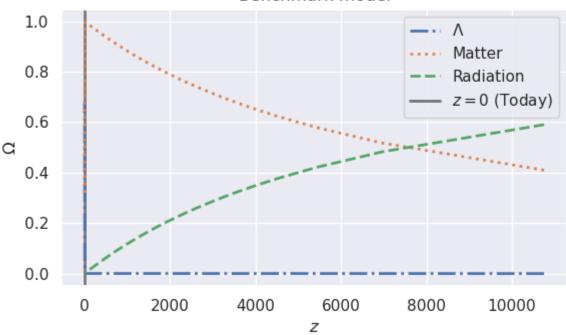
ax.plot(df_past['z'], df_past['dark density'], label = '$\Lambda$', linestyle = 'dashdo
ax.plot(df_past['z'], df_past['matter density'], label = 'Matter', linestyle = 'dotted'
ax.plot(df_past['z'], df_past['radiation density'], label = 'Radiation', linestyle = 'd

ax.axvline(x = 0, color = 'black', linestyle = 'solid', linewidth = 2, alpha = 0.5, lab
ax.set_xlabel('$z$')
ax.set_ylabel('$\Omega$')
ax.set_ylabel('$\Omega$')
ax.set_title('Evolution of densities as a function of $z>0$.\nBenchmark model')

plt.legend(loc = 'best')

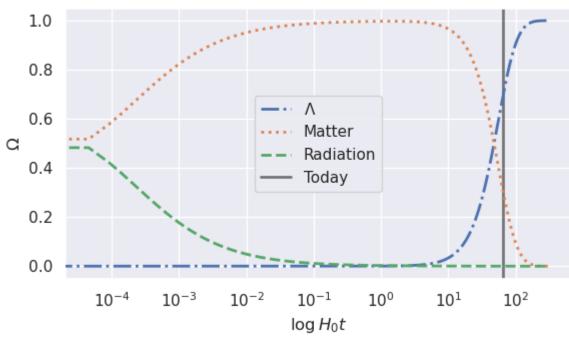
plt.tight_layout()
plt.savefig('./figs/densities_evolution_z_future.png', dpi = 300)
plt.show()
plt.close(fig)
```

Evolution of densities as a function of z > 0. Benchmark model



```
In [18]:
          # Densities evolution as a function of time (log scale)
          fig, ax = plt.subplots(figsize = (6, 4))
          ax.plot((df['t']-min(df['t']))*H_0, df['dark density'], label = '$\Lambda$', linestyle
          ax.plot((df['t']-min(df['t']))*H 0, df['matter density'], label = 'Matter', linestyle =
          ax.plot((df['t']-min(df['t']))*H_0, df['radiation density'], label = 'Radiation', lines
          ax.axvline(x = (-min(df['t']))*H_0, color = 'black', linestyle = 'solid', linewidth = 2
          ax.set xlabel('$\log{H Ot}$')
          ax.set ylabel('$\Omega$')
          ax.set_title('Evolution of densities as a function of time.\nBenchmark model')
          # X-scale to exponential
          plt.xscale('log')
          plt.legend(loc = 'best')
          plt.tight_layout()
          plt.savefig('./figs/densities evolution logt.png', dpi = 300)
          plt.show()
          plt.close(fig)
```

Evolution of densities as a function of time. Benchmark model



```
In [19]: # The time of the universe as predicted by the benchmark model
# APROXIMATION
time = (-min(df['t']))*H_0 # Units: km/s/Mpc

# Remove km and Mpc from time
time = 1/(time*(3.2478e-20))/31536000/1e9 # Units: Gyr
print('Time in Gyr: ', time)
```

Time in Gyr: 15.191131561951128

Point 3

Compute the proper, the luminosity and angular-diameter distance. Make plots of such quantities for the base cosmology, and for variations of it. You can get some inspiration by the plots in chapter 5 and 6 of the Barbara Ryden book[2].

We are going to use the benchmark model to calculate the proper, the luminosity and angular-diameter distance, and we will plot the distance as a function of redshift. For our benchmark model, with k=0, we have:

Proper distance.

$$egin{aligned} d_p(t_0) &= c \int_{t_e}^{t_0} rac{dt}{a(t)} \quad ; \quad H = rac{\dot{a}}{a} \implies rac{da}{dt} = aH \implies dt = rac{da}{aH} \ \implies d_p &= c \int_a^1 rac{1}{a} rac{da}{aH} = c \int_a^1 rac{da}{a\dot{a}} \end{aligned}$$

Luminosity distance.

$$d_L=d_p(1+z)$$

Angular diameter distance.

```
d_A=rac{d_L}{(1+z)^2}
```

```
In [20]: # Filter only positive redshifts of benchmark model
    df_distance = df[df['z'] > 0.0]
    df_distance.head(5)
```

```
Out[20]:
                                 matter density radiation density
                                                                  dark density
                                                                                      a dot
             -0.954140 0.000093
                                       0.409850
                                                         0.590150 7.294826e-13 6086.261121 10754.290258
             -0.954139 0.000144
                                       0.517432
                                                         0.482568 3.389521e-12 4359.339198
                                                                                              6965.110766
              -0.954138 0.000183
                                       0.577814
                                                         0.422186 7.871165e-12 3651.396884
                                                                                              5456.598931
              -0.954138 0.000218
                                       0.619015
                                                         0.380985 1.410864e-11 3237.775637
                                                                                              4596.172906
             -0.954137 0.000248
                                       0.649805
                                                         0.350195 2.205987e-11 2957.104080
                                                                                              4024.427866
```

```
In [21]:

def proper_distance(df):
    """
    Function to calculate the proper distance of a given simulation.
    INPUT:
        - df: DataFrame of a simulation
    OUTPUT:
        - dp: List wit proper distance of the simulation
    """

# List for proper distance
d_p = [0]
        n = len(df)

# Iterate over values in dataframe (a, a_dot)
for i in range(1, len(df)):
        delta_a = (df['a'][n-i]-df['a'][n-i-1]) # Psudo-differential
        d_p.append(d_p[-1] + delta_a/(df['a'][n-i]*df['a_dot'][n-i])) # Proper distance

# Return list
    return np.array(d_p[::-1])*H_0 # Adjust units to Hubble distance
```

In [22]:
Get proper distance of three simulations (dark, matter, benchmark)
df_distance['d_p'] = proper_distance(df_distance)

/tmp/ipykernel_543/3701085558.py:2: SettingWithCopyWarning:
A value is trying to be set on a copy of a slice from a DataFrame.
Try using .loc[row_indexer,col_indexer] = value instead

See the caveats in the documentation: https://pandas.pydata.org/pandas-docs/stable/user_guide/indexing.html#returning-a-view-versus-a-copy df_distance['d_p'] = proper_distance(df_distance)

```
In [23]:
```

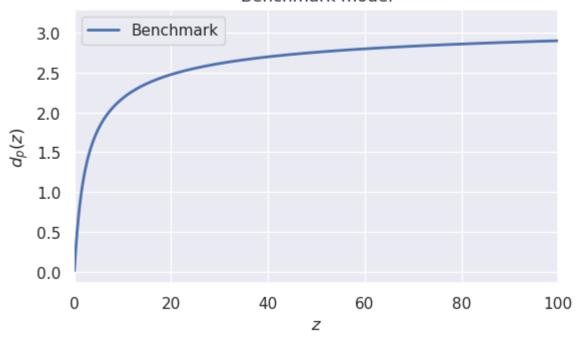
```
# Filter out high redshifts
df_distance = df_distance[(df_distance['z'] > 0.01) & (df_distance['z'] < 1000)]</pre>
```

```
In [24]:
# Proper distance evolution as a function of redshift
fig, ax = plt.subplots(figsize = (6, 4))

ax.plot(df_distance['z'], df_distance['d_p'], label = 'Benchmark', linestyle = 'solid',
    ax.set_xlabel('$z$')
    ax.set_ylabel('$d_p(z)$')
    ax.set_title('Proper distance as a function of $z$.\nBenchmark model')
    ax.set_xlim(0, 100)
    ax.legend()

plt.tight_layout()
    plt.savefig('./figs/proper_distance_z.png', dpi = 300)
    plt.show()
    plt.close(fig)
```

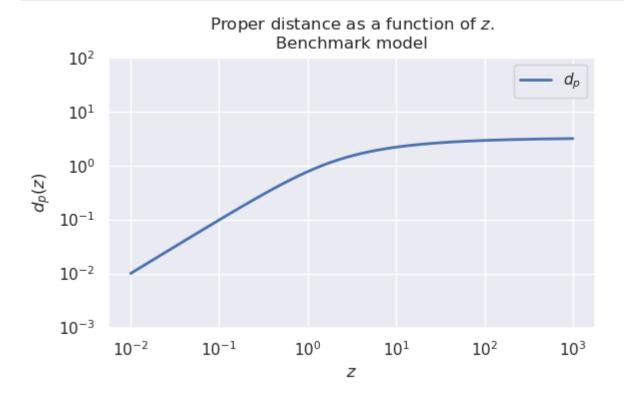
Proper distance as a function of z. Benchmark model



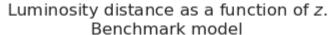
```
In [25]: # Proper distance evolution as a function of redshift (log-log scale)
fig, ax = plt.subplots(figsize = (6, 4))
ax.plot(df_distance['z'], df_distance['d_p'], label = '$d_p$', linestyle = 'solid', lin
ax.set_xlabel('$z$')
ax.set_ylabel('$d_p(z)$')
ax.set_title('Proper distance as a function of $z$.\nBenchmark model')
plt.xscale('log')
plt.yscale('log')
ax.set_ylim(0.001, 100)
plt.legend()

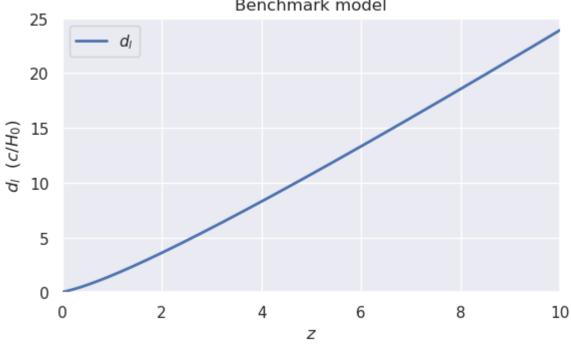
plt.tight_layout()
plt.savefig('./figs/proper_distance_logz.png', dpi = 300)
```

```
plt.show()
plt.close(fig)
```



```
In [26]:
          df_distance['d_1'] = df_distance['d_p']*(1+df_distance['z']) # Compute Luminosity dista
          df_distance['d_a'] = df_distance['d_l']/(1+df_distance['z'])**2 # Compute angular diame
In [27]:
          # Luminosity distance evolution as a function of redshift
          fig, ax = plt.subplots(figsize = (6, 4))
          ax.plot(df_distance['z'], df_distance['d_1'], label = '$d_1$', linestyle = 'solid', lin
          ax.set xlabel('$z$')
          ax.set ylabel('$d 1$ $(c/H 0)$')
          ax.set_title('Luminosity distance as a function of $z$.\nBenchmark model')
          ax.set_xlim(0, 10)
          ax.set_ylim(0, 25)
          plt.legend()
          plt.tight_layout()
          plt.savefig('./figs/luminosity_distance_z.png', dpi = 300)
          plt.show()
          plt.close(fig)
```





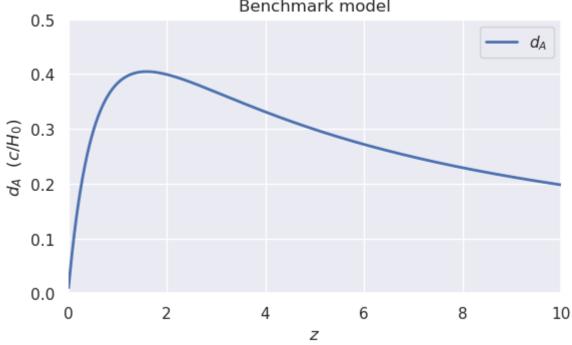
```
In [28]: # Angular diameter distance evolution as a function of redshift
fig, ax = plt.subplots(figsize = (6, 4))

ax.plot(df_distance['z'], df_distance['d_a'], label = '$d_A$', linestyle = 'solid', lin

ax.set_xlabel('$z$')
ax.set_ylabel('$d_A$ $(c/H_0)$')
ax.set_title('Angular-diameter distance as a function of $z$.\nBenchmark model')
ax.set_xlim(0, 10)
ax.set_ylim(0, 0.5)
plt.legend()

plt.tight_layout()
plt.savefig('./figs/angular_distance_z.png', dpi = 300)
plt.show()
plt.close(fig)
```





```
In [29]:
          # Plot of distances as a function of redshift
          fig, ax = plt.subplots(figsize = (6, 4))
          ax.plot(df_distance['z'], df_distance['d_a'], label = '$d_A$', linestyle = 'solid', lin
          ax.plot(df distance['z'], df distance['d l'], label = '$d l$', linestyle = 'solid', lin
          ax.plot(df_distance['z'], df_distance['d_p'], label = '$d_p$', linestyle = 'solid', lin
          ax.set xlabel('$z$')
          ax.set_ylabel('$d$')
          ax.set_title('Different distances as a function of $z$.\nBenchmark model')
          ax.set xlim(0, 10)
          ax.set_ylim(0, 3)
          plt.legend()
          plt.tight_layout()
          plt.savefig('./figs/distances_z.png', dpi = 300)
          plt.show()
          plt.close(fig)
```



Point 4.

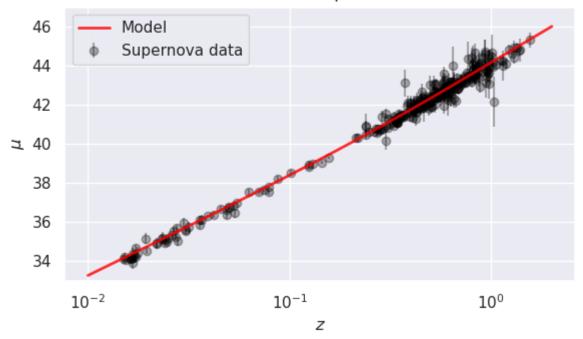
In particular use the Luminosity distance to compute the distance modulus

$$\mu = 5log_{10}rac{d_L}{1Mpc} + 25$$

and compare it against the data provided.

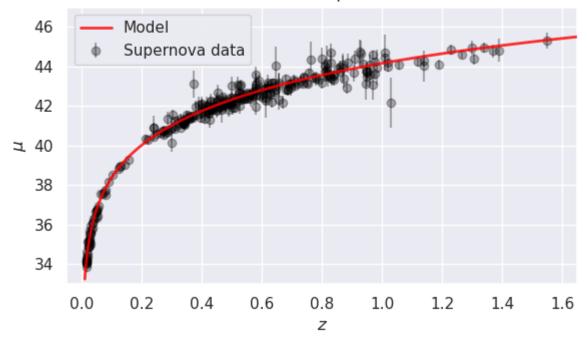
```
In [30]:
          df_distance['mu'] = 5*np.log10(df_distance['d_1']*c/H_0) + 25 # Compute distance modulu
In [31]:
          # Load and clean up supernova data
          import requests
          response = requests.get('https://supernova.lbl.gov/Union/figures/SCPUnion mu vs z.txt')
          data = response.text
          data = data.split('\n')[4:]
          data = [float(value) for line in data for value in line.split('\t')[1:]]
          data = np.array(data)
          data = data.reshape(len(data)//3, 3)
In [32]:
          df supernova = pd.DataFrame(data, columns = ['z', 'mu', 'mu err']) # Dataframe of super
          df_aux = df_distance[(df_distance['z'] > 0.00) & (df_distance['z'] < 2)] # Filter out h</pre>
In [33]:
          # Plot of distance modulus as a function of redshift (model vs. supernova data, log sca
          fig, ax = plt.subplots(figsize = (6, 4))
          ax.errorbar(df supernova['z'], df supernova['mu'], df supernova['mu err'],
```

Distance modulus as a function of z. Simulation vs Supernova data



```
plt.tight_layout()
plt.savefig('./figs/supernova_linear.png', dpi = 300)
plt.show()
plt.close(fig)
```

Distance modulus as a function of z. Simulation vs Supernova data



In []: