

Homework 2

Aerosol Physics, Chemistry, Clouds and Climate

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Problem 1 Condensation

1.a

Assumed values:

- $T = 298.15K$
- $P = 1atm$
- $C_{\infty} = 5E7 \text{ molecules } cm^{-3} = 8.14E(-12) \text{ kg } m^{-3}$
- $C_s = 0$
- $D_g = 1E(-5) \text{ m}^2 \text{ s}^{-1}$
- Equation used for the condensation diameter growth rate:

$$\frac{dD_p}{dt} = \frac{4D_g}{\rho D_p} (C_{\infty} - C_s) \beta$$

The figure 1.a shows the condensation diameter growth rate vs diameter. The condensation growth rate is independent of the diameter of the particle in the kinetic regime ($D_p < 10nm$). Whereas in the continuum regime ($D_p > 1000nm$), the condensation diameter growth rate is inversely proportional to the diameter of the particle. In the transition regime ($10nm < D_p < 1000nm$), the condensation diameter growth rate is a function of the diameter of the particle and behaves similar to a sigmoid function between the kinetic and continuum regimes.

1.b

In order to calculate condensation mass growth rate the following equation was used:

$$\frac{dM_P}{dt} = J = 2\pi D_g D_p (C_{\infty} - C_s) \beta$$

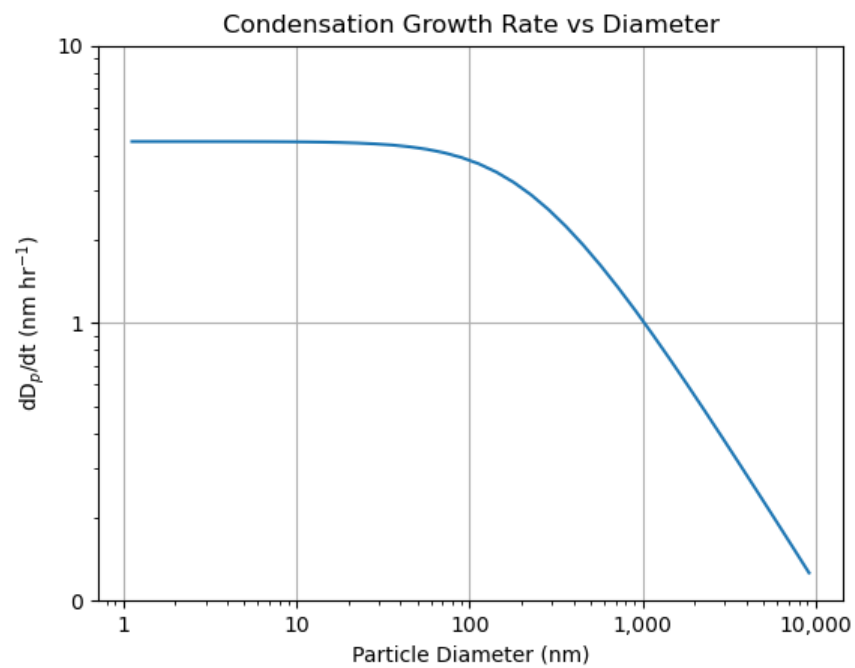


Figure 1: Problem(1) part(a) Condensation diameter growth rate vs diameter

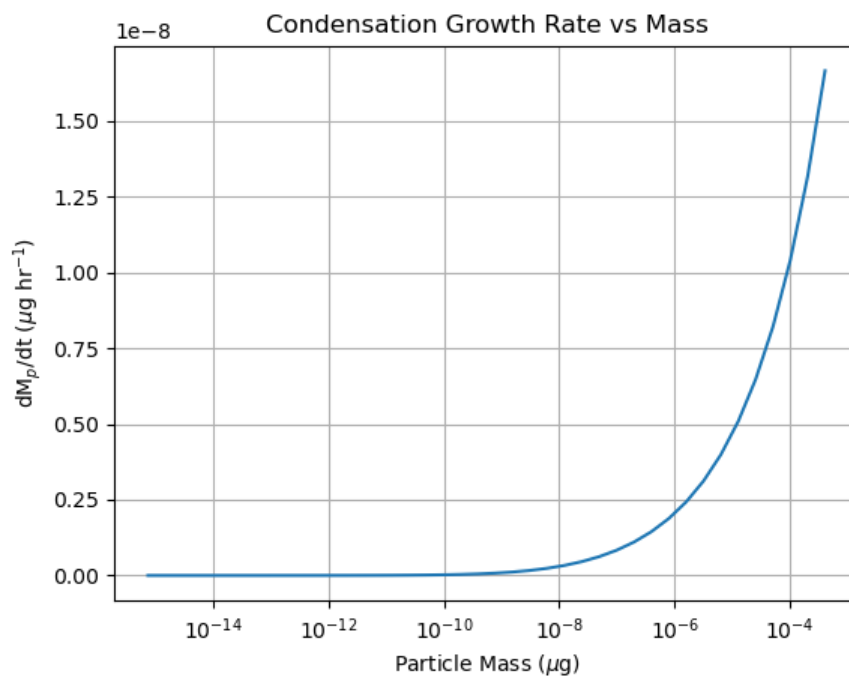


Figure 2: Problem(1) part(b) Condensation mass growth rate vs particle diameter

Problem 2 Coagulation

2.a Coagulation kernel and rate

Assumed values:

- $T = 298.15K$
- $P = 1atm$
- $N_1 = 3000 \text{ cm}^{-3}$, $D_{p1} = 10nm$
- $N_2 = 400 \text{ cm}^{-3}$, $D_{p2} = 100nm$
- $D_g = 1E(-5) \text{ m}^2 \text{ s}^{-1}$

Fuchs form of the brownian coagulation coefficient K_{12} :

$$K_{12} = 2\pi(D_1 + D_2)(D_{p1} + D_{p2})\left(\frac{D_{p1} + D_{p2}}{D_{p1} + D_{p2} + 2(g_1^2 + g_2^2)^{1/2}} + \frac{8(D_1 + D_2)}{(c_1^2 + c_2^2)^{1/2}(D_{p1} + D_{p2})}\right)^{-1}$$

$$K_{12} = 2.38E(-8) \text{ cm}^3 \text{ s}^{-1}$$

$$K_{table13.3} = 2.5E(-8) \text{ cm}^3 \text{ s}^{-1}$$

$$J_{12} = K_{12}N_1N_2 = 2.85E(-2) \text{ cm}^{-3} \text{ s}^{-1}$$

2.b Frequency of collision

The frequency of collision with the smaller particles = $K_{12}N_1 = 7.13E(-5)$

$$\text{The frequency of collision} = \frac{\text{Collision rate}}{N_1}$$

2.c Growth rate due to coagulation

The growth rate due to coagulation is given by the following equation:

$$\frac{dV_{p2}}{dt} = k_{12}N_1V_{p1}$$

$$\frac{dD_{p2}}{dt} = k_{12}N_1V_{p1}\frac{2}{\pi D_{p2}^2}$$

$$\frac{dD_{p2}}{dt} = 8.56E(-3)nm \text{ hr}^{-1}$$

2.d Growth in 1 week

The percent change in diameter in 1 week = 1.44%

The percent change in mass in 1 week = 4.31%

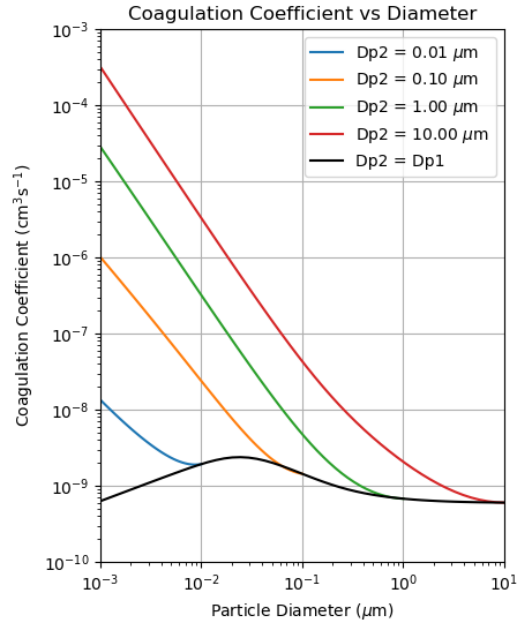


Figure 3: Fuchs Coagulation the code

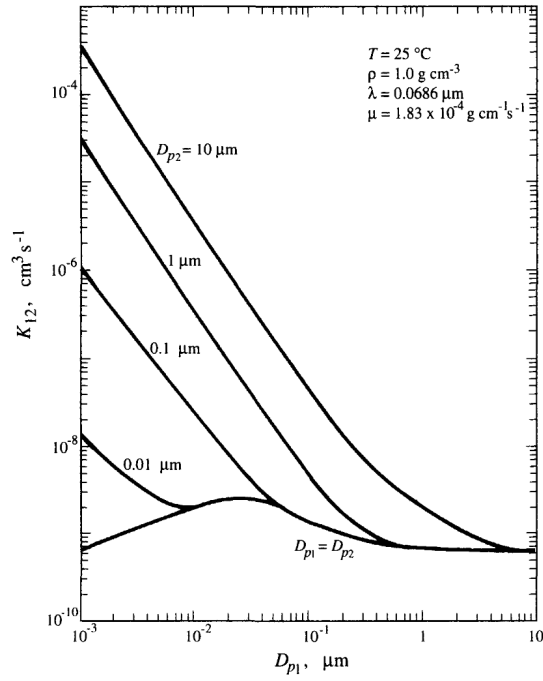


Figure 4: Table_13.5 Fuchs Coagulation plot

HW_2_Aero_cond_coag

February 21, 2024

1 Condensation and Coagulation

Homework 2 for the course Aerosols physics and chemistry

2 Libraries

```
[33]: #importing libraries
import numpy as np
import matplotlib.pyplot as plt
import matplotlib as mpl
from pyfluids import Fluid, FluidsList, Input
from scipy import constants as const
import pandas as pd

mpl.rcParams["figure.dpi"] = 100
```

3 Global constants

```
[34]: h = const.Planck # J s Planck constant
r_gas_c = const.R # J/(mol K) molar gas constant
k_boltz = const.Boltzmann # J/K Boltzmann constant
g = const.g # m/s2 acceleration due to gravity
n_avo = const.Avogadro # 1/mol Avogadro's number
```

4 Function Definitions

```
[35]: # Default air standard properties
pressure_std = 101325 # Pa
temperature_std = 273.15 + 25 # K #??????? need to make sure this is the right_
↪ temperature
air_std = Fluid(FluidsList.Air).with_state(Input.pressure(pressure_std), Input.
↪ temperature(temperature_std-273.15))

def c_cunningham(d_p, lamda = 65E-9):
    #defining the function
    #Cunningham correction factor
```

```

#Dp is the particle diameter in meters
#lamda is the mean free path of the gas in meters
#C is the Cunningham correction factor
kn = 2 * lamda / d_p
return 1 + kn * (1.257 + 0.4 * np.exp(-1.1 / kn))

def mean_free_path(temperature, pressure):
# mean free path of air calculator
#T is the temperature in Kelvin
#P is the pressure in Pascals
#lamda is the mean free path of the gas in meters
R = 8.314 # J/(mol K) gas constant
M = 0.0289647 # kg/mol molar mass of air
air = Fluid(FluidsList.Air).with_state(Input.pressure(pressure), Input.
temperature(temperature - 273.15))
viscosity = air.dynamic_viscosity # Pa s dynamic viscosity
return 2 * viscosity / (pressure * np.sqrt(8 * M / (np.pi * R *
temperature)))

def reynolds_number(d_p, velocity, fluid_density = air_std.density,
dynamic_viscosity = air_std.dynamic_viscosity):
#Reynolds number calculator
#Dp is the particle diameter in meters
#rho_f is the density of the fluid in kg/m^3
#g is the acceleration due to gravity in m/s^2
#C is the Cunningham correction factor
#Re is the Reynolds number
return (d_p * fluid_density * velocity) / dynamic_viscosity

def settling_velocity(d_p, rho_p, temperature, pressure):
# Check if Dp_input is an array or a single value
if np.isscalar(d_p):
d_p_array = np.array([d_p]) # Convert to array for uniform processing
else:
d_p_array = d_p # Use the array as is

velocities = [] # Empty list to store calculated velocities
for d_p in d_p_array: # Process each Dp individually
g = 9.81 # m/s^2
l_mfp = mean_free_path(temperature, pressure)
c_cun = c_cunningham(d_p, l_mfp)
air = Fluid(FluidsList.Air).with_state(Input.pressure(pressure), Input.
temperature(temperature - 273.15))
mu_f = air.dynamic_viscosity

```

```

    rho_f = air.density
    s_velocity = c_cun * (rho_p * g * d_p**2) / (18 * mu_f) # Stokes
    ↪settling velocity
    Re = reynolds_number(d_p, s_velocity, fluid_density=rho_f,
    ↪dynamic_viscosity=mu_f)
    if Re < 1:
        velocities.append(s_velocity)
    else:
        # Adjusted iterative approach for Re > 1, similar to before
        m_p = np.pi * rho_p * d_p**3 / 6
        for i in range(100):
            # c_d = 24 / Re * (1 + 0.15 * Re**(0.687)) # Updated drag
            ↪coefficient expression
            c_d = 24 / Re * (1 + 3/16 * 0.43 * Re)
            # s_velocity = np.sqrt((4 * m_p * g) / (3 * np.pi * c_d * rho_f
            ↪* Dp**2))
            s_velocity = np.sqrt((m_p * g) / (1/8 * np.pi * c_d * rho_f *
            ↪d_p**2))
            Re_new = reynolds_number(d_p, s_velocity, fluid_density=rho_f,
            ↪dynamic_viscosity=mu_f)
            if abs(Re_new - Re) < 0.01:
                break # Exit the loop if the change in Reynolds number is
            ↪small enough
            else:
                Re = Re_new
                velocities.append(s_velocity)

    velocities_array = np.array(velocities) # Convert list to array

    if np.isscalar(d_p):
        return velocities_array[0] # Return a single value if input was scalar
    else:
        return velocities_array # Return array if input was array

def cond_dia_growth_rate(d_p, c_inf, c_sat, diff_coefficient, rho_p = 1E3):
    #condensation diameter growth rate
    #Dp is the particle diameter in meters
    b_factor = beta_corr_cond(d_p)
    d_dp_dt = b_factor * 4 * diff_coefficient * (c_inf - c_sat) / (rho_p * d_p)
    return d_dp_dt
    # # if np.isscalar(d_p):
    # #     d_p_array = np.array([d_p]) # Convert to array for uniform
    ↪processing
    # # else:
    # #     d_p_array = d_p # Use the array as is
    # # dia_growth_rate = [] # Empty list to store calculated velocities

```



```

#
# return dia_growth_rate

def cond_mass_growth_rate(d_p, c_inf, c_sat, diff_coefficient_vapor, rho_p = 1E3):
    #condensation mass growth rate
    #Dp is the particle diameter in meters
    #c_inf is the concentration of the condensing vapor in the bulk gas in
    #molec/m^3
    #c_sat is the saturation concentration of the condensing vapor in molec/m^3
    #diff_coefficient is the diffusion coefficient of the condensing vapor in
    #m^2/s
    #rho_p is the density of the particle in kg/m^3 (default is 1E3 kg/m^3)
    #m_p is the mass of the particle in kg
    #particle mass
    # J is the same as condensation mass growth rate
    j_coag_calculated = 2 * np.pi * diff_coefficient_vapor * d_p * (c_inf -
    c_sat) * beta_corr_cond(d_p) #kg/s
    return j_coag_calculated

def beta_corr_cond(d_p, lamda = 65E-9):
    #Dahneke correction factor
    #Dp is the particle diameter in meters
    #B is the Darkener correction factor
    # kn is the Knudsen number
    # lamda is the mean free path of the gas in meters
    kn = 2 * lamda / d_p
    b_factor = (1 + kn) / (1 + 2 * kn * (1 + kn))
    return b_factor

def unit_conversion(conversion, variable=1):
    switcher = {
        #length
        'cm3_to_m3': variable * 1E-6,
        'm3_to_cm3': variable * 1E6,
        #time
        'hr_to_s': variable * 3600,
        's_to_hr': variable * 1/3600,
        #volume
        'm3_to_L': variable * 1000,
        'L_to_m3': variable * 1/1000,
        #mass
        'kg_to_g': variable * 1000,
        'g_to_kg': variable * 1/1000,
    }

```

```

        'kg_to_ug': variable * 1E9,
        'ug_to_kg': variable * 1E-9,

        #pressure
        'Pa_to_kPa': variable * 1E-3,
        'kPa_to_Pa': variable * 1E3,
        'atm_to_Pa': variable * 101325,
        'Pa_to_atm': variable * 1/101325,
        #temperature
        'C_to_K': variable + 273.15,
        'K_to_C': variable - 273.15,
    }
    return switcher.get(conversion, 'Invalid conversion')

def dp_to_mp_distribution(d_p, rho_p = 1E3):
    #Dp is the particle diameter in meters
    #rho_p is the density of the particle in kg/m^3
    #m_p is the mass of the particle in kg
    m_p = np.pi * rho_p * d_p**3 / 6
    return m_p

def vol_sphere(d_p):
    #Dp is the particle diameter in meters
    #V is the volume of the sphere in m^3
    return (4/3) * np.pi * (d_p/2)**3

def surf_sphere(d_p):
    #Dp is the particle diameter in meters
    #A is the surface area of the sphere in m^2
    return 4 * np.pi * (d_p/2)**2

```

5 Other Functions - Coagulation

```

[36]: def k_coag(d_p_1, d_p_2, temp = temperature_std, rho_1 = 1000, rho_2 = 1000): #
        # Source: Sienfeld, J. H., & Pandis, S. N. (2006). Atmospheric chemistry_
        ↪and physics: from air pollution to climate change. John Wiley & Sons.
        # 2nd Edition Table 13.1
        #coagulation coefficient
        #Dp1 is the diameter of particle 1 in meters
        #Dp2 is the diameter of particle 2 in meters
        #c1 is the concentration of particle 1 in molec/m^3
        #c2 is the concentration of particle 2 in molec/m^3

```

```

#d_diff1 is the diffusion coefficient of particle 1 in m^2/s
#d_diff2 is the diffusion coefficient of particle 2 in m^2/s
#temp is the temperature in Kelvin
#rho_1 is the density of particle 1 in kg/m^3
#rho_2 is the density of particle 2 in kg/m^3
# k_bolt is the Boltzmann constant in J/K defined in the global constants

# Convert dp_1 and dp_2 to numpy arrays if they are not already
dp_1_np = np.array([d_p_1])
dp_2_np = np.array([d_p_2])

#k is the coagulation coefficient
# used the mass of a particle because the c_1 and c_2 are velocity of the
particles
air_01 = Fluid(FluidsList.Air).with_state(Input.pressure(pressure_std),
Input.temperature(temp-273.15))
vis_air = air_01.dynamic_viscosity # Pa s dynamic viscosity
d_diff1 = k_boltz * temp * np.array([c_cunningham(dp) for dp in dp_1_np]) /
(3 * np.pi * vis_air * dp_1_np)
d_diff2 = k_boltz * temp * np.array([c_cunningham(dp) for dp in dp_2_np]) /
(3 * np.pi * vis_air * dp_2_np)
m_1 = rho_1 * np.pi * dp_1_np ** 3 / 6 # mass of one particle in kg
m_2 = rho_2 * np.pi * dp_2_np ** 3 / 6 # mass of one particle in kg
c_1 = np.sqrt(8 * k_boltz * temp / (np.pi * m_1))
c_2 = np.sqrt(8 * k_boltz * temp / (np.pi * m_2))
l_1 = 8 * d_diff1 / (np.pi * c_1)
l_2 = 8 * d_diff2 / (np.pi * c_1)
g_1 = 1 / (3 * dp_1_np * l_1) * ((dp_1_np + l_1) ** 3 - (dp_1_np ** 2 + l_1
** 2) ** (3 / 2)) - dp_1_np
g_2 = 1 / (3 * dp_2_np * l_2) * ((dp_2_np + l_2) ** 3 - (dp_2_np ** 2 + l_2
** 2) ** (3 / 2)) - dp_2_np
denominator = (((dp_1_np + dp_2_np) / (dp_1_np+dp_2_np+2*np.
sqrt(g_1**2+g_2**2))) + 8*(d_diff1+ d_diff2)/(c_1**2+c_2**2)**0.5 /
(dp_1_np+dp_2_np))

k_coag_fuchs = 2 * np.pi * (dp_1_np + dp_2_np) * (d_diff1 + d_diff2) /
denominator
return k_coag_fuchs

```

6 Bins definition

```

[37]: #defining the bins
bin_number = 40
bins_lower = np.geomspace(1e-9, 10.3e-6, bin_number + 1) #
bins_upper = bins_lower[1:]
bins_lower = bins_lower[:-1]

```

```
bins_g_mean = np.sqrt(bins_lower * bins_upper) # geometric mean
bins_g_mean_mass = dp_to_mp_distribution(bins_g_mean)
```

7 Problem 1

Condensation of Sulfuric Acid (H₂SO₄)

```
[38]: molec_weight_H2SO4 = 98.079 / 1000 #kg/mol
c_inf_h2so4 = 5E7 / const.centimeter * 3 / const.Avogadro * molec_weight_H2SO4 #unit
    ↪ conversion from molec/cm^3 kg/m^3
c_sat_h2so4 = 0 #molec/m^3
diff_h2so4 = 1E-5 #m^2/s
```

8 Problem 1-part a

Condensation diameter growth rate

```
[39]: dp_dt_condensation = cond_dia_growth_rate(bins_g_mean, c_inf_h2so4,
    ↪ c_sat_h2so4, diff_h2so4, rho_p = 1E3)
dm_dt_condensation = cond_mass_growth_rate(bins_g_mean, c_inf_h2so4,
    ↪ c_sat_h2so4, diff_h2so4, rho_p = 1E3)
```

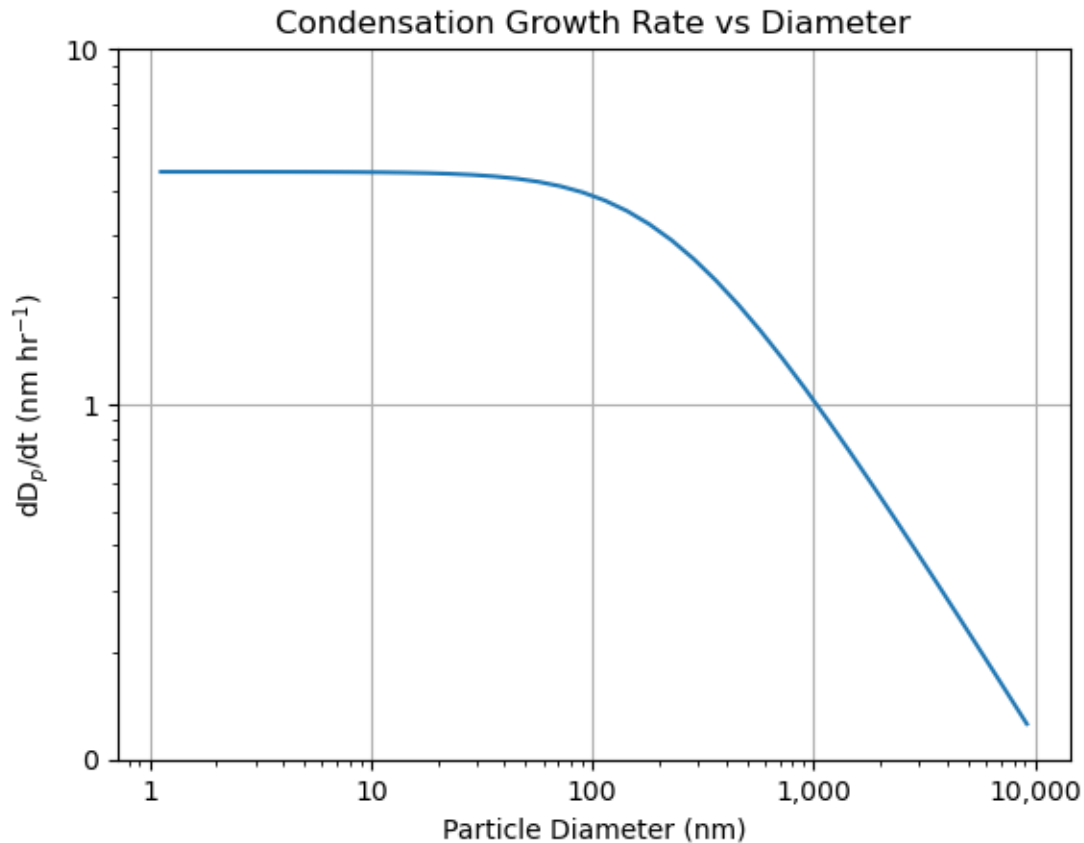
```
[40]: # plot condensation growth rate vs diameter
fig, ax = plt.subplots()
ax.plot(bins_g_mean / const.nano, dp_dt_condensation * 1e9 * 3600)
ax.set_xscale('log')
ax.set_yscale('log')
ax.set_xlabel('Particle Diameter (nm)')
ax.set_ylabel(r'dD$_p$/dt (nm hr$^{-1}$)')
ax.set_title('Condensation Growth Rate vs Diameter')
ax.set_ylim([1e-1, 1e1])

# change the x-axis labels to float
formatter = mpl.ticker.FuncFormatter(lambda x, pos: '{:,.0f}'.format(x))
ax.xaxis.set_major_formatter(formatter)

# change the y-axis labels to float
formatter = mpl.ticker.FuncFormatter(lambda x, pos: '{:,.0f}'.format(x))
ax.yaxis.set_major_formatter(formatter)

plt.grid()

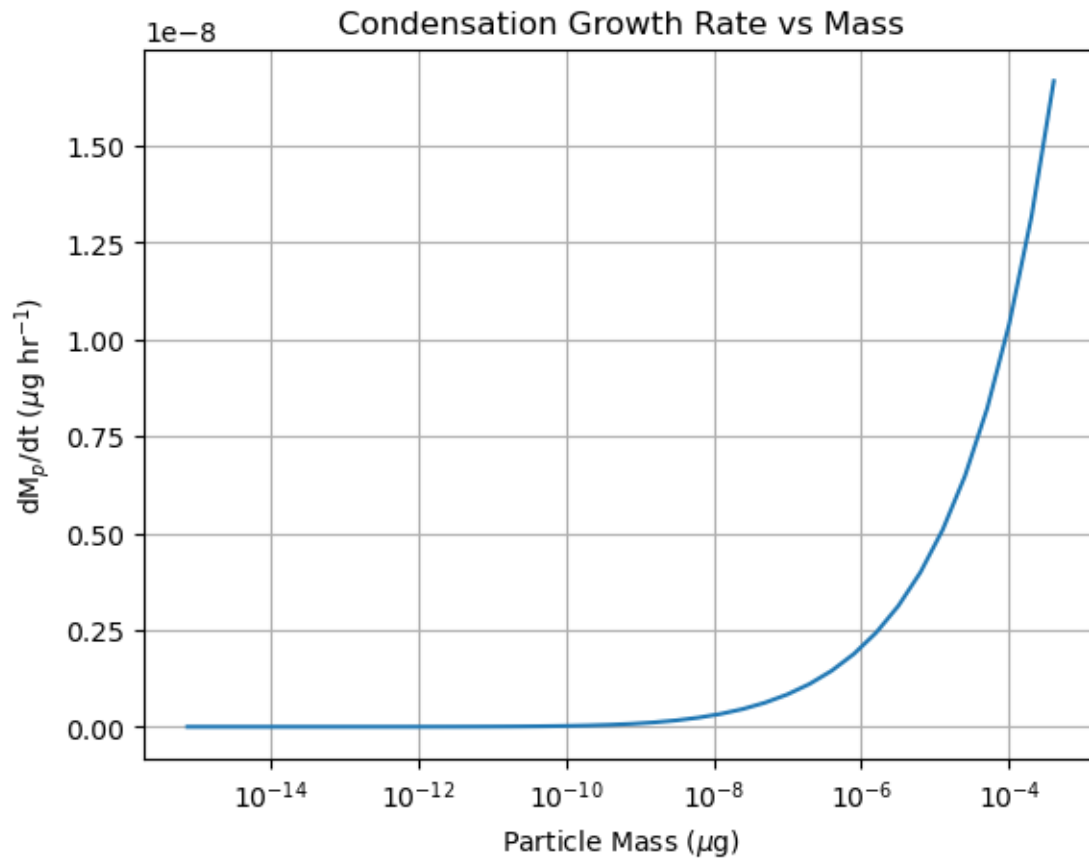
fig.savefig('hw2_pr1_1_condensation_growth_rate_vs_diameter.png')
plt.show()
```



9 Problem 1-part b

Condensation mass growth rate

```
[41]: # plot mass growth rate vs particle mass
fig, ax = plt.subplots()
ax.plot(bins_g_mean_mass * unit_conversion('kg_to_ug'),
        dm_dt_condensation * unit_conversion('kg_to_ug') *
        unit_conversion('hr_to_s')) #converted to kg/s to ug/hr
#converted kg to ug
ax.set_xscale('log')
ax.set_xlabel('Particle Mass ( $\mu\text{g}$ )')
ax.set_ylabel('dMp/dt ( $\mu\text{g hr}^{-1}$ )')
ax.set_title('Condensation Growth Rate vs Mass')
plt.grid()
plt.savefig('hw2_pr1_2_condensation_growth_rate_vs_mass.png')
plt.show()
```



10 Problem 2

Coagulation as a growth mechanism

```
[42]: # plot coagulation coefficient vs particle diameter
fig, ax = plt.subplots(figsize=(5, 6))

for dp_2 in np.geomspace(1e-8, 1e-5, 4):
    dp_1 = np.geomspace(1e-9, dp_2, 50)
    k_c = k_coag(dp_1, dp_2).flatten() # Flatten the array to make it 1D
    ax.plot(dp_1 / const.micro, k_c / const.centimeter ** 3, label=f'Dp2 = {dp_2 / const.micro:.2f} $\mu\text{m}$')

dp_same = np.geomspace(1e-9, 1e-5, 50)
k_c = k_coag(dp_same, dp_same).flatten()
ax.plot(dp_same / const.micro, k_c / const.centimeter ** 3, label=f'Dp2 = Dp1', color='black')

ax.set_xscale('log')
```

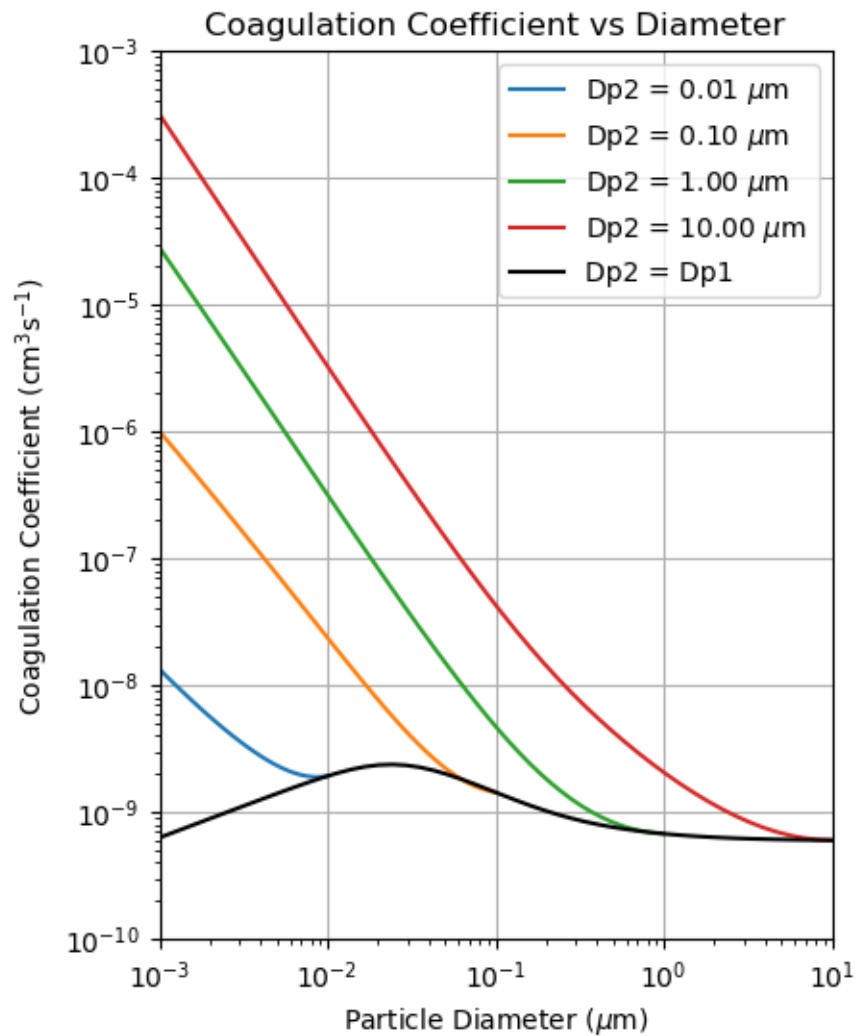
```

ax.set_yscale('log')
ax.set_xlabel(f'Particle Diameter ( $\mu\text{m}$ )')
ax.set_ylabel(r'Coagulation Coefficient ( $\text{cm}^3\text{s}^{-1}$ )')
plt.grid()
ax.set_title('Coagulation Coefficient vs Diameter')
ax.set_xlim([1e-3, 10]) # Set the limits of x-axis
ax.set_ylim([1e-10, 1e-3]) # Set the limits of y-axis
ax.legend()

# Adjust the space on the left side of the plot
plt.subplots_adjust(left=0.2)

plt.savefig('hw2_pr2_1_coagulation_coefficient_vs_diameter.png')
plt.show()

```



```
[43]: # making a table of the coagulation coefficient
dp_2 = [2e-9, 1e-8, 1e-7, 1e-6, 1e-5, 2e-5]
dp_1 = [1e-9, 1e-8, 1e-7, 1e-6, 1e-5, 2e-5]
k_c = np.array([[k_coag(d1, d2) for d1 in dp_1] for d2 in dp_2])
# print in the table format pandas dp2 as the index and dp1 as the columns
k_c_squeezed = np.squeeze(k_c, axis=-1) # Squeeze the array along the last axis
df = pd.DataFrame(k_c_squeezed / const.centimeter**3, index = dp_2, columns = dp_1)
```

```
[44]: N1 = 3000 / const.centimeter**3 #particle/m^3
dp1 = 10 * const.nanometer #m
N2 = 400 / const.centimeter**3 #particle/m^3
dp2 = 100 * const.nanometer #m
rho_1 = rho_2 = 1E3 #kg/m^3
j_coag = k_coag(dp1, dp2) * N1 * N2 #collision/m^3/s
print(f'{k_coag(dp1, dp2)[0] / const.centimeter**3:.2e} cm^3/s') # Convert
    ↳coagulation coefficient from m^3/s to cm^3/s
print(f'{j_coag[0] * const.centimeter**3:.2e} collision/cm^3/s') # Convert
    ↳collision/m^3 /s to collision/cm^3 s
# converted m^3/s to cm^3/s
```

2.38e-08 cm³/s

2.85e-02 collision/cm³/s

```
[45]: # frequency of collisions with particles
freq_coll = k_coag(dp1, dp2)[0] * N1 #collision/s
print(f' frequency of collisions with smaller particles: {freq_coll:.2e}
    ↳collision/s')

#delete this
# one_N2_particles_surface = np.pi * dp2**2 / 4 #m^2/particle
# all_N2_particles_surface = one_N2_particles_surface * N2 #m^2/m^3
# freq_coll = j_coag * all_N2_particles_surface
```

frequency of collisions with smaller particles: 7.13e-05 collision/s

```
[46]: # growth rate of the larger particles
volume_of_dp1 = vol_sphere(dp1) #m^3
mass_growth_rate_dp2_coag = freq_coll * volume_of_dp1 #m^3/s
dia_growth_rate_dp2_coag = mass_growth_rate_dp2_coag * (2/ np.pi / dp2**2) #m/s
dia_growth_rate_dp2_coag_nm_hr = dia_growth_rate_dp2_coag / const.nanometer * 3600
    ↳#nm/hr
mass_growth_rate_dp2_coag_kg_hr = mass_growth_rate_dp2_coag * 3600 #kg/hr
print(f'Growth rate of the larger particles: {dia_growth_rate_dp2_coag_nm_hr:.
    ↳2e} nm/hr')
```

Growth rate of the larger particles: 8.56e-03 nm/hr


```
[47]: # Total growth after 1 week
total_growth_dia = dia_growth_rate_dp2_coag_nm_hr * 24 * 7 #nm
percent_growth_dia = total_growth_dia / (dp2 / const.nano) * 100
total_growth_mass = mass_growth_rate_dp2_coag_kg_hr * 24 * 7 #kg
percent_growth_mass = total_growth_mass / (vol_sphere(dp2)) * 100

print(f'Total growth in diameter after 1 week: {total_growth_dia:.2f} nm')
print(f'Percent growth in diameter after 1 week: {percent_growth_dia:.2f} %')
print(f'Total growth in mass after 1 week: {total_growth_mass:.2e} kg')
print(f'Percent growth in mass after 1 week: {percent_growth_mass:.4f} %')
```

```
Total growth in diameter after 1 week: 1.44 nm
Percent growth in diameter after 1 week: 1.44 %
Total growth in mass after 1 week: 2.26e-23 kg
Percent growth in mass after 1 week: 4.3122 %
```