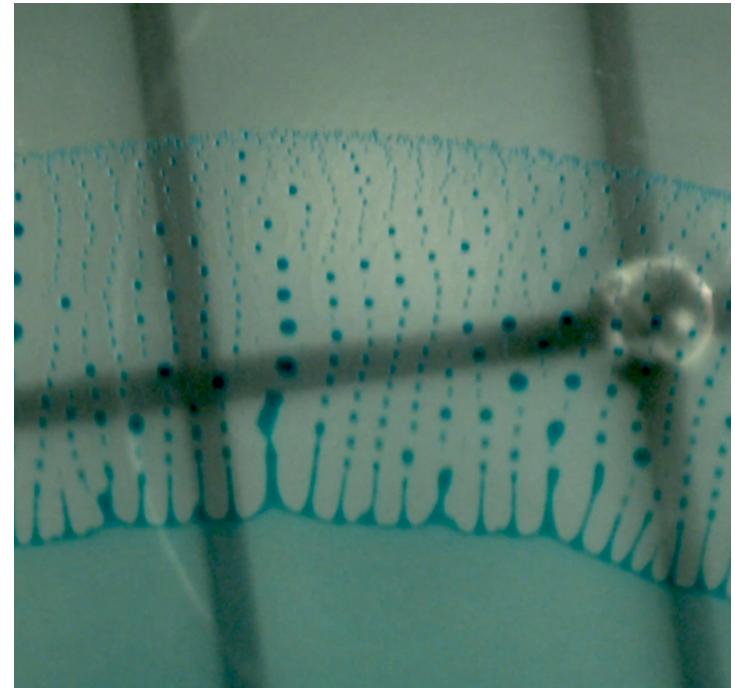


10. Droplet Explosion

Yuezhen (Lily) Dong

*"When a drop of a water mixture (e.g. water-alcohol) is deposited on the surface of a hydrophobic liquid (e.g. vegetable oil), the resulting drop may sometimes **fragment** into smaller droplets. Investigate the parameters that affect the **fragmentation** and the size of the final droplets."*

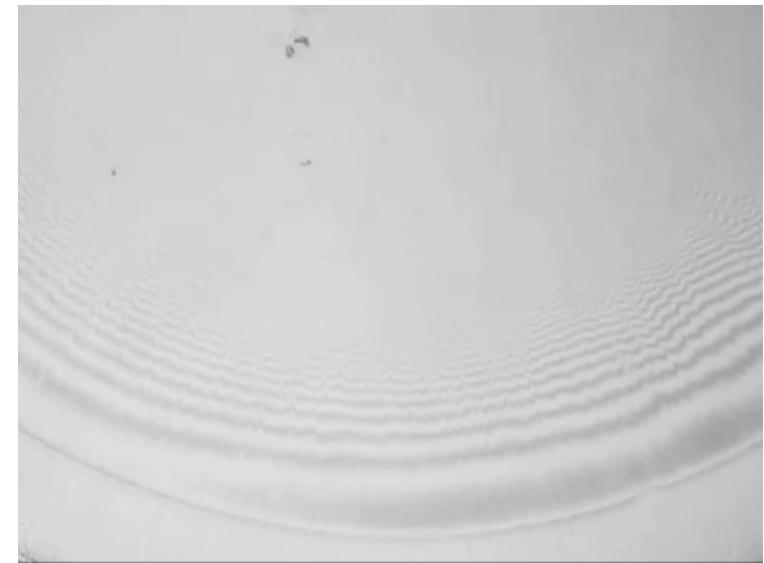


Problem Statement

"When a drop of a water mixture (e.g. water-alcohol) is deposited on the surface of a hydrophobic liquid (e.g. vegetable oil), the resulting drop may sometimes fragment into smaller droplets. Investigate the parameters that affect the fragmentation and the size of the final droplets."

Parameters:

1. *Initial concentration of alcohol*
2. *Initial volume of droplets*
3. *Depth of the hydrophobic liquid*
4. *Types of solvent*



Overview

1

Phenomenon

Reproduction and Explanation of the Phenomenon

2

Experimental Setup

Measurement Techniques, Camera Views

3

Theoretical Model

Qualitative Explanation, Quantitative Models

4

Key Parameter Interactions

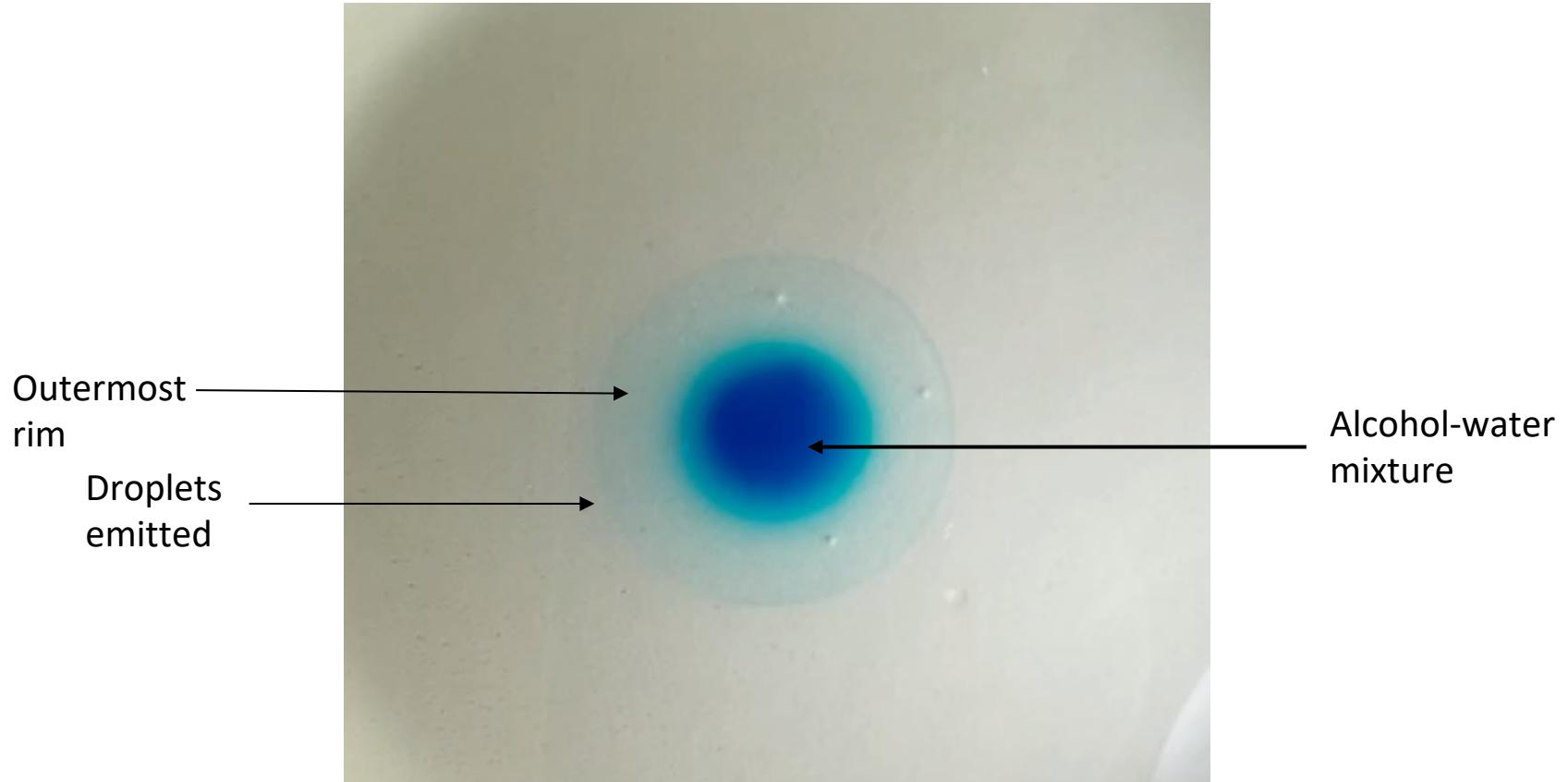
Effects of Changing Physical Parameters

5

Conclusion

Further Insights and General Investigations

Phenomenon



Experimental Setup

Introduction

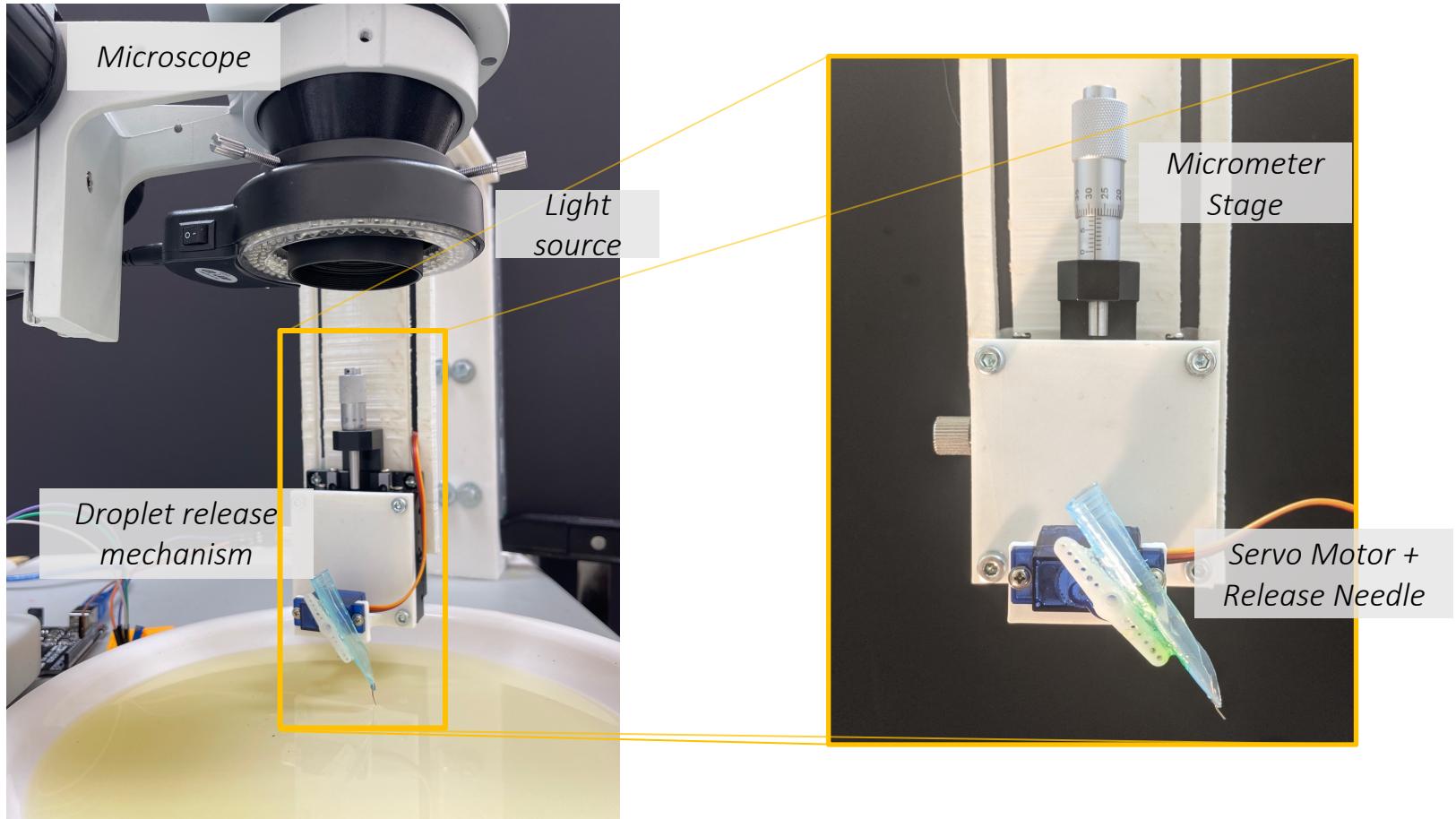
Experimental Setup

Theoretical Model

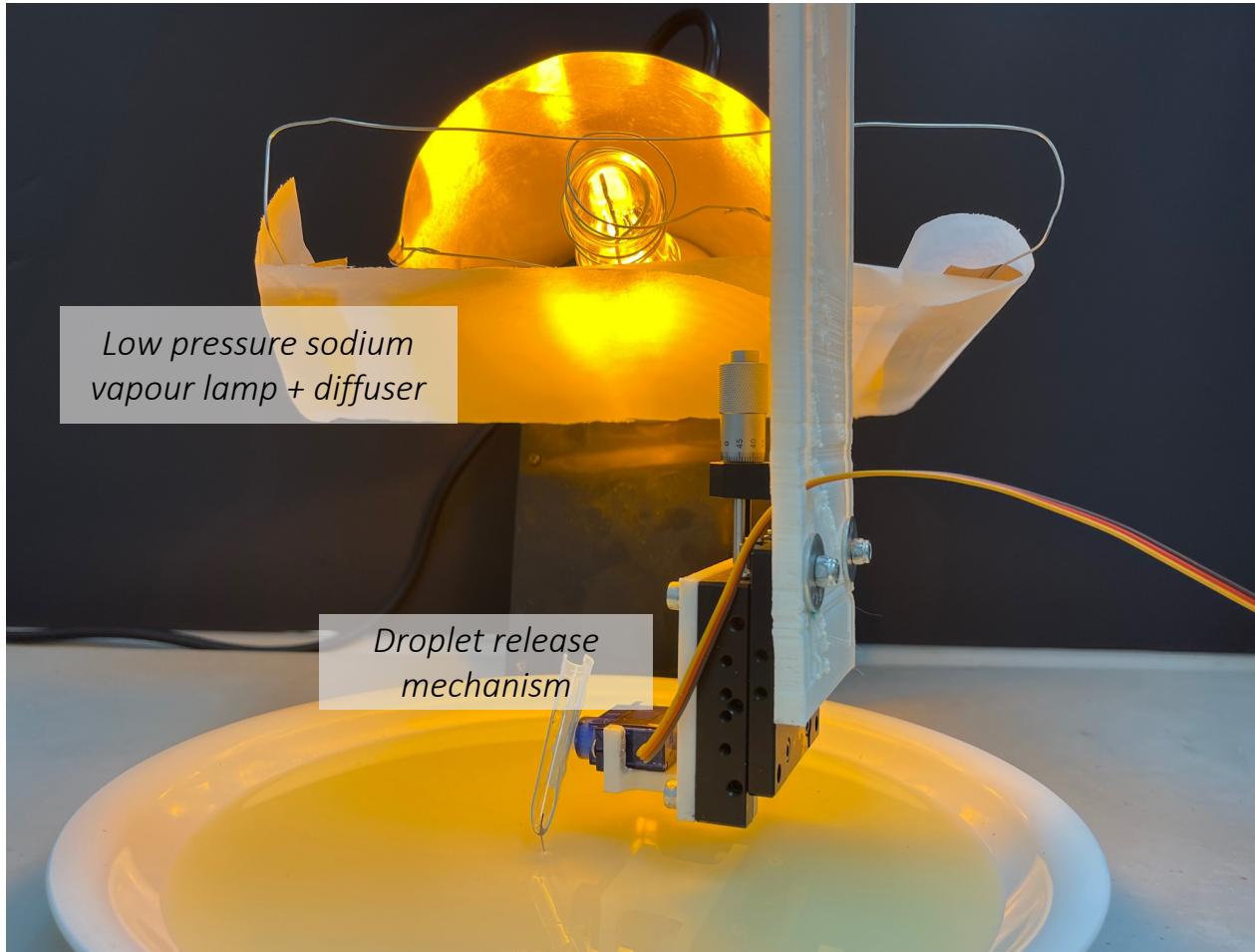
Key Parameters

Conclusion

Microscopy Setup



Interferometry Setup



Introduction

Experimental Setup

Theoretical Model

Key Parameters

Conclusion

Experimental Measurement



Syringes ($\pm 0.5mL$)



Micropipette
($\pm 0.6\%$)



200 mL graduated cylinder
($\pm 2mL$)

Theoretical Model

Introduction

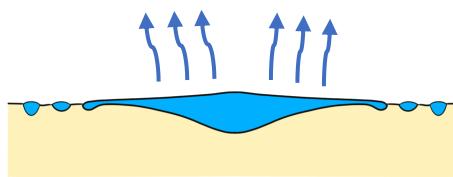
Experimental Setup

Theoretical Model

Key Parameters

Conclusion

Theoretical Model



Alcohol
Evaporation

Interfacial Tension
& Shear Stress

Droplet
Fragmentation

Geometry

j_v = evaporation rate of alcohol

d = width of the rim

h_m = minimum thickness of the droplet expansion

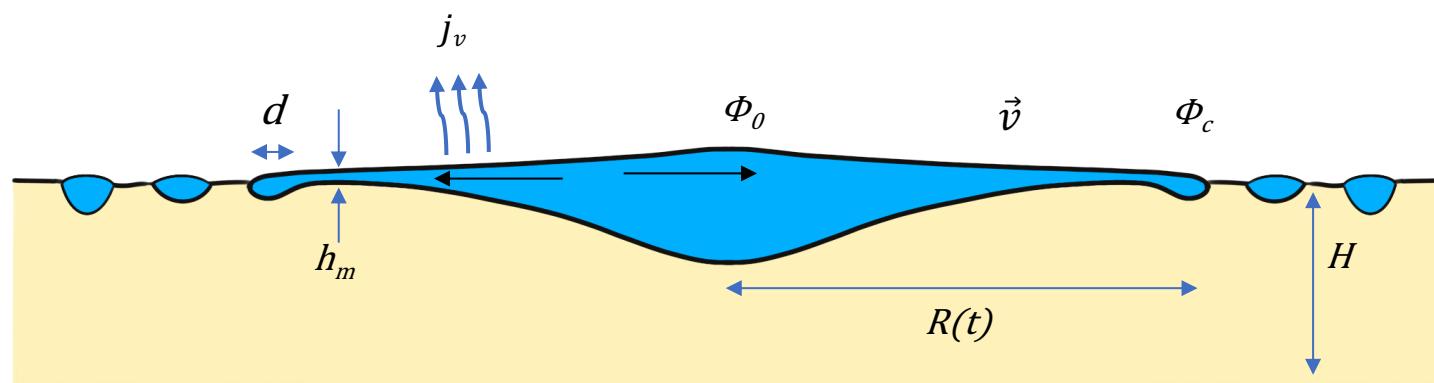
Φ_0 = initial mass fraction of alcohol in the center of the drop

Φ_c = mass fraction of alcohol at the rim of the expansion

\vec{v} = velocity of the Marangoni flow

$R(t)$ = radius of droplet

H = depth of the oil layer



Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Assumptions

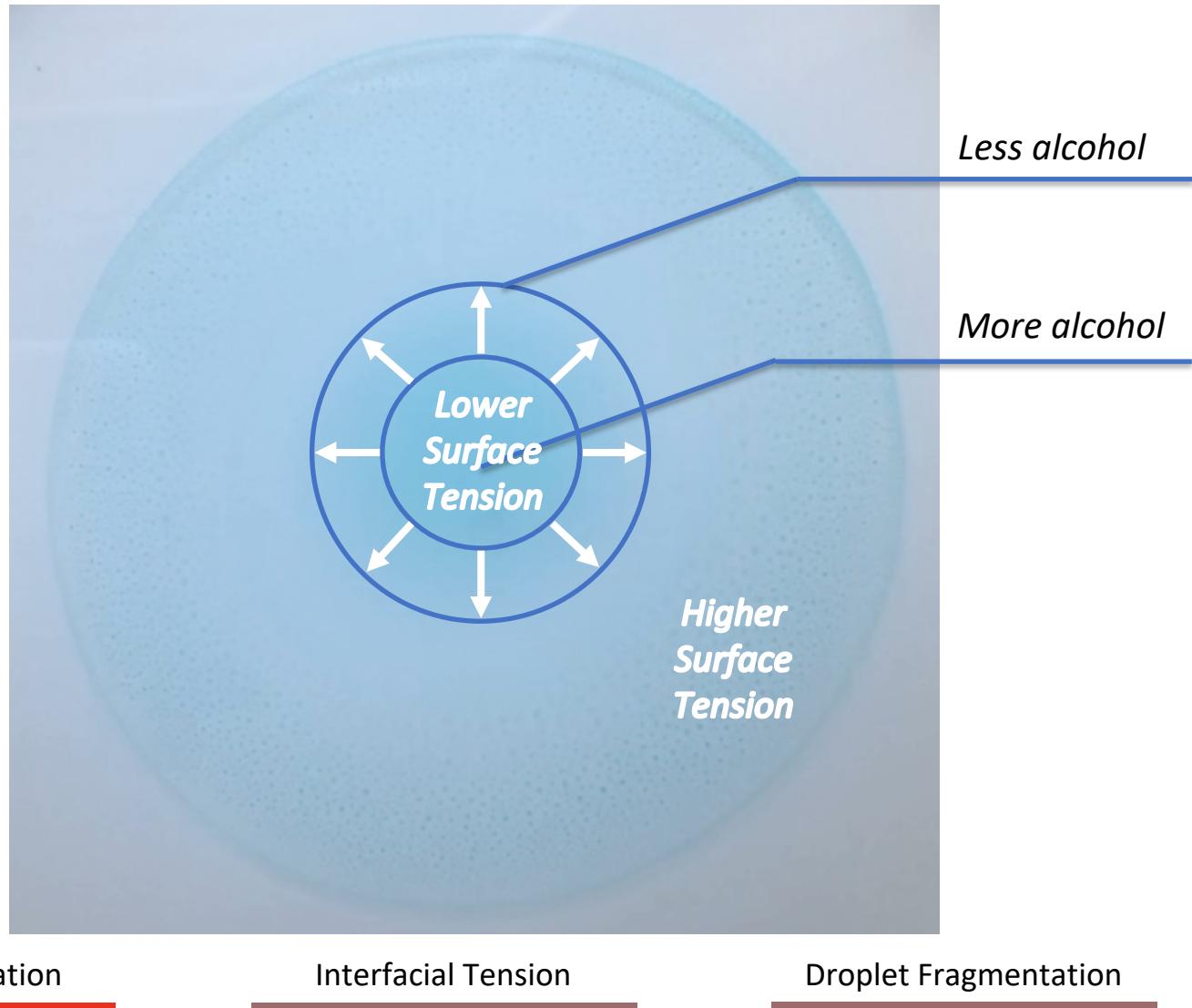
- *The thinner film of the droplet is caused by alcohol evaporation*
- *The evaporation of water is negligible*
- *Ambient room temperature remains constant*
- *The impact of water based dye is negligible*

Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Marangoni Flow



Evaporation Flux [8]

To calculate the evaporation flux, use Fick's law,

$$j_m = -D \nabla c \cdot \mathbf{n}$$

j_m is the evaporation flux

\mathbf{n} is the unit normal vector

c is the molar concentration of IPA

Considering the mass conservation across the interface, this leads to

$$\rho_l (\mathbf{u}_l \cdot \mathbf{n} - u_\Gamma) = j_m$$

u_Γ is the normal velocity of the interface

\mathbf{n} is the unit normal vector

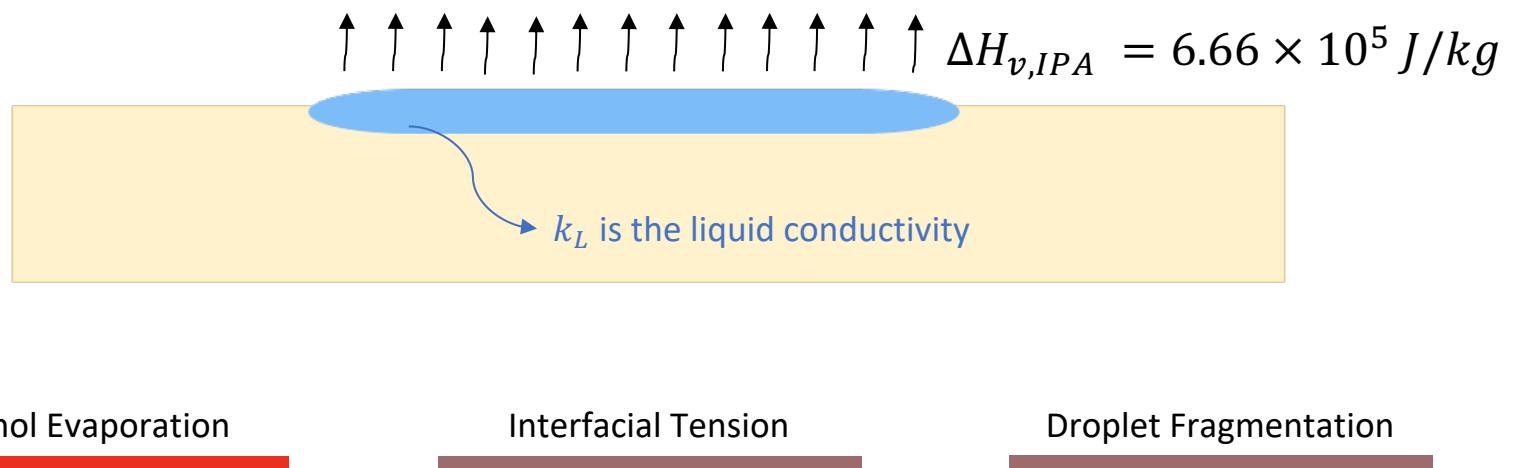
c is the molar concentration of IPA

Interfacial Energy Balance

For negligible heat conduction and convection in the air, the interfacial energy balance is,

$$-k_L \mathbf{n} \cdot \nabla T = \Delta H_v j(x)$$

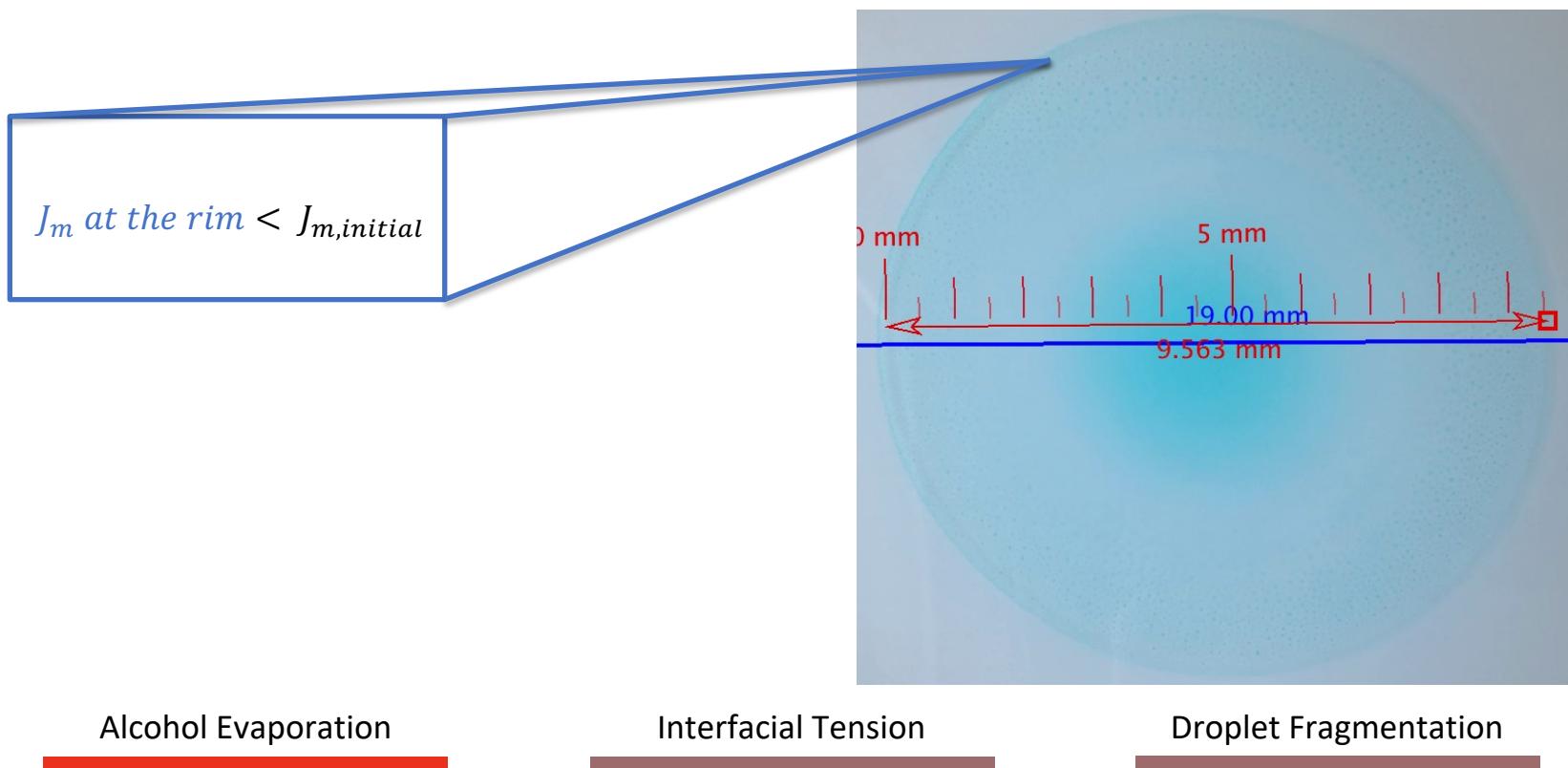
ΔH_v is the specific latent heat of evaporation, IPA = $6.66 \times 10^5 \text{ J/kg}$



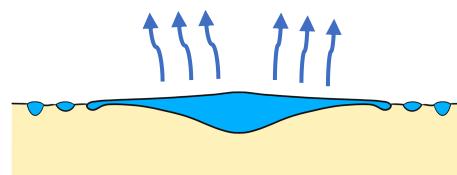
Alcohol Evaporation

The spreading parameter, $S = \gamma_{oa} - \gamma_{ma} - \gamma_{mo}$

Evaporation and total Mass flux, $J_m = \int j_m d\Gamma$



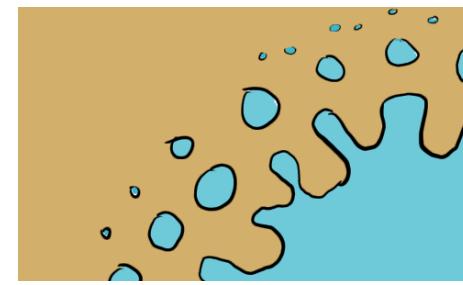
Theoretical Model



Alcohol
Evaporation

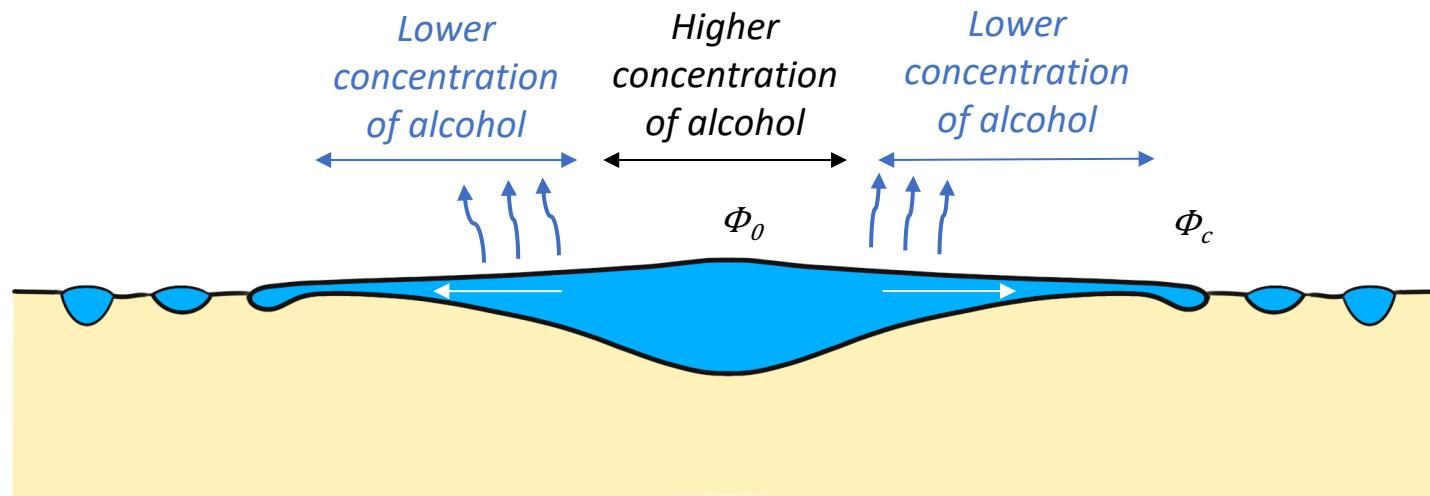


Interfacial Tension
& Shear Stress



Droplet
Fragmentation

Geometry



Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Assumptions

1. *Ambient room temperature is constant.*
2. *The dropping height remains constant throughout the experiment.*
3. *The dimensions of the oil plate remains constant.*
4. *The effects caused by the ingredients of water-based food dye are negligible.*
5. *The alcohol evaporation during the dropping process is negligible.*
6. *The evaporation of water in this experiment is negligible.*

Alcohol Evaporation

Interfacial Tension

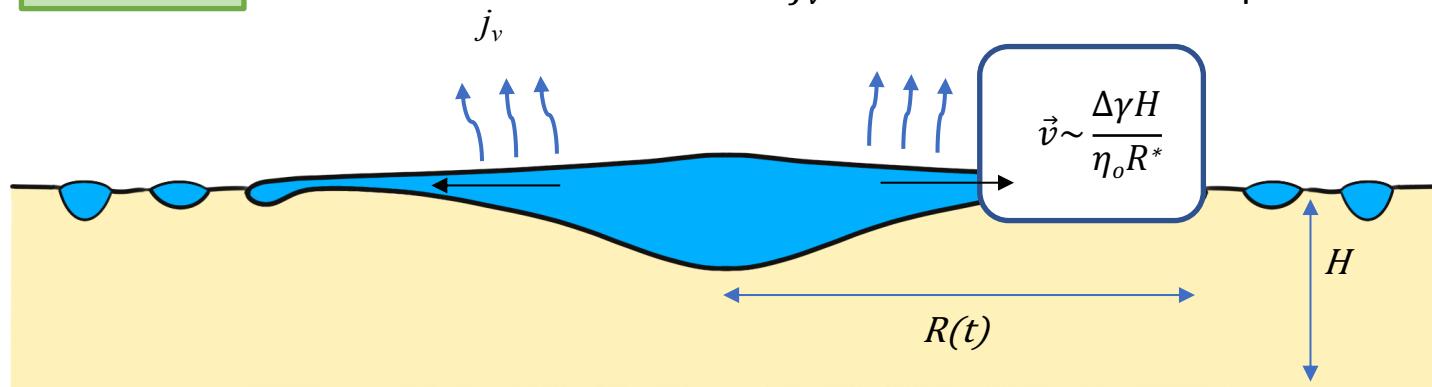
Droplet Fragmentation

Marangoni Flow

Surface tension gradient is the same magnitude of $\frac{\Delta\gamma}{R^*}$

$$\tau \sim \frac{R^*}{V}$$

Time scale of the experiment



Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

$$\Delta\gamma = \gamma_c - \gamma_o$$

γ_c = interfacial tension corresponding to Φ_c

γ_o = interfacial tension corresponding to Φ_0

R^* = characteristic radius of the drop

η_o = viscosity of oil

j_v = volume of alcohol evaporated

Marangoni Flow [7]

The evolution equation for the height, h , as a function of radial coordinate r and time t ,

$$\frac{\partial h}{\partial t} = \frac{1}{r\mu} \frac{\partial}{\partial r} \left(\frac{rh^3}{3} \frac{\partial p}{\partial r} - \frac{rh^2}{2} \frac{\partial \sigma}{\partial r} \right) + \omega_e$$

Pressure Gradient
Surface Tension Gradient

p is the pressure in the droplet

ω_e is the evaporative volume flux

μ is viscosity of droplet

σ is surface tension

Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Marangoni Flow

The pressure in the droplet is dominated by surface tension. Therefore, the pressure p can be given by the Laplace pressure:

$$p = -\frac{1}{r} \frac{\partial}{\partial r} \left(\sigma \frac{r \partial_r h}{\sqrt{1 + (\partial_r h)^2}} \right)$$

The boundary and initial conditions of $h(r,t)$ are,

$$\left(\frac{\partial h}{\partial r} \right)_{r=0} = 0$$

$$h(r,t) = 0$$

$$h(r,0) = h_0 r$$

Marangoni Flow [1]

Solved for characteristic radius of the droplet expansion

$$R^* \sim \left(\frac{(\Phi_0 - \Phi_c) \Delta \gamma H \Omega_0}{(1 - \Phi_c) \eta_o j_v} \right)^{\frac{1}{4}}$$

$$\tau \sim \left(\frac{(\Phi_0 - \Phi_c) \eta_o \Omega_0}{(1 - \Phi_0) \Delta r H j_v} \right)^{\frac{1}{2}}$$

Solved time scale based on the convection of the liquid from the center to the edge

$$(0.4 < \Phi_0 < 0.8) \\ 1\text{mm} < H < 20\text{ mm} \\ 2\text{mm}^3 < \Omega_0 < 500\text{ mm}^3)$$

Restrictions applied

$$R_{max} \sim 0.28 (\pm 0.02) R^* \\ t_{exp} \sim 1.5 (\pm 0.15) \tau$$

Estimated maximum radius and correlated time

[1] Keiser, L., Bense, H., Colinet, P., Bico, J., Reyssat, E., 2017.

Alcohol Evaporation

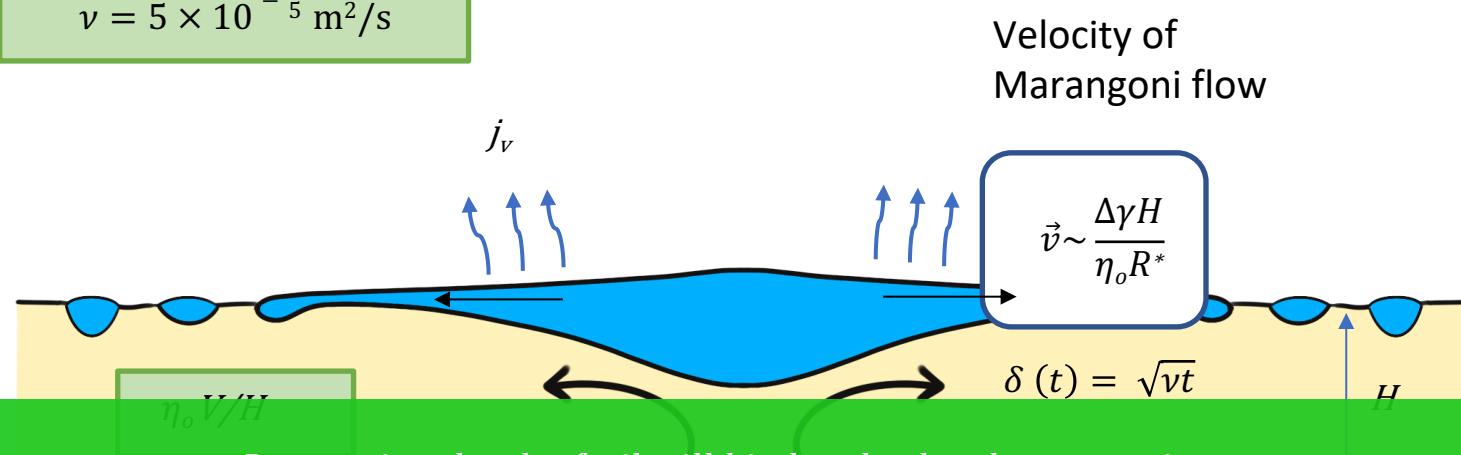
Interfacial Tension

Droplet Fragmentation

Viscous stress

Kinematic viscosity of oil

$$\nu = 5 \times 10^{-5} \text{ m}^2/\text{s}$$



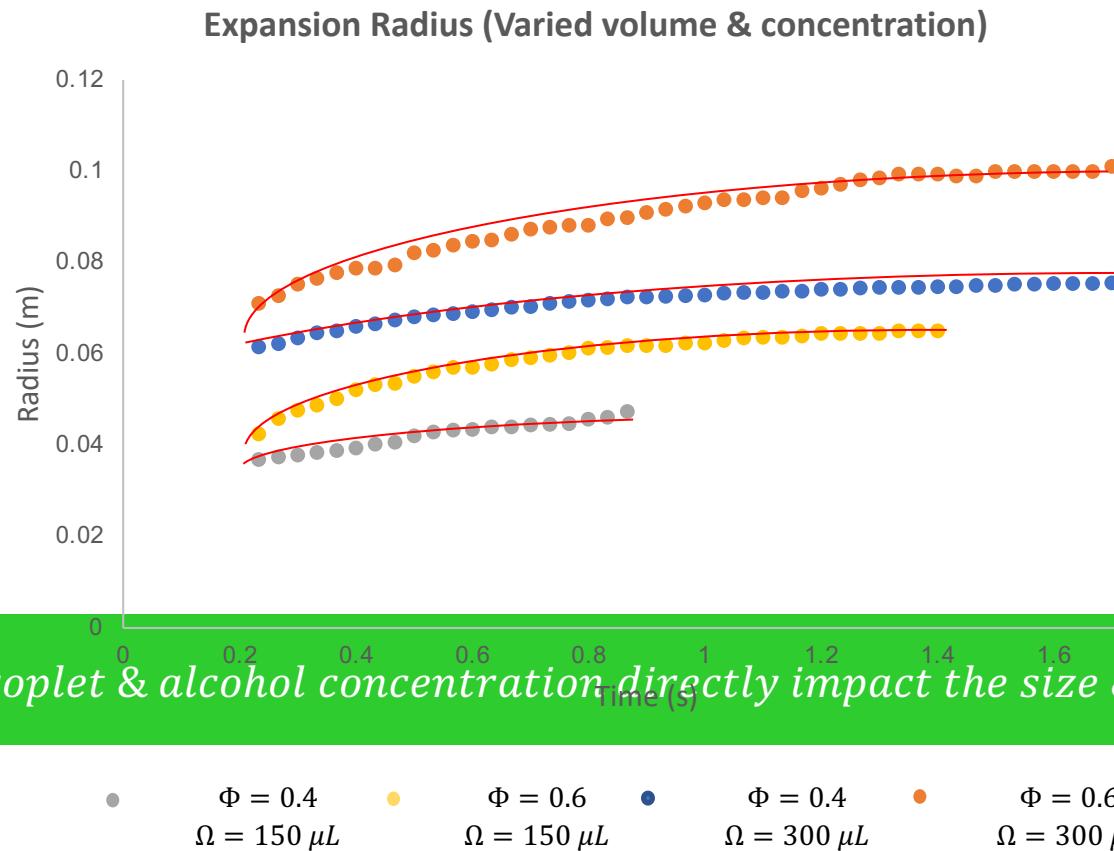
Decreasing depth of oil will hinder the droplet expansion
 Order of magnitude
 of the viscous stress

Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Preliminary Experiment Results

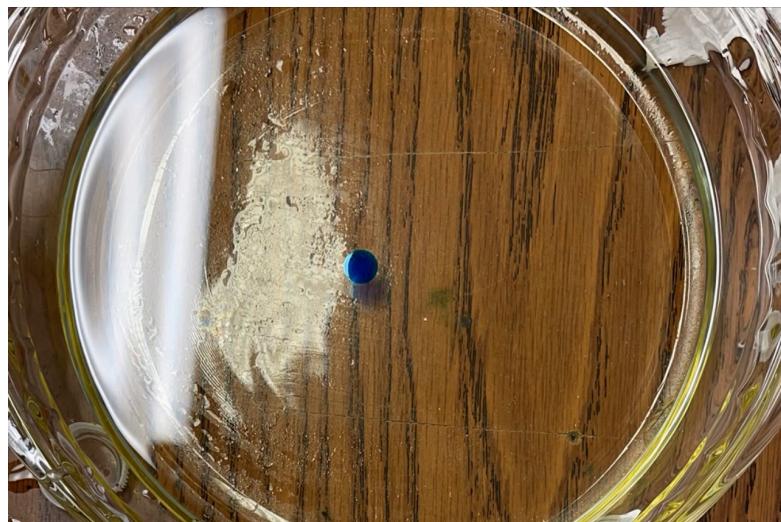


Alcohol Evaporation

Interfacial Tension

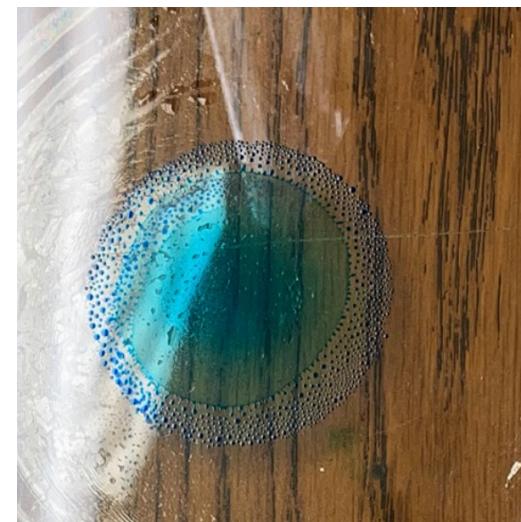
Droplet Fragmentation

Critical Point of the Spreading Parameter - Ethanol



$$\Phi = 0.5$$

Alcohol Evaporation

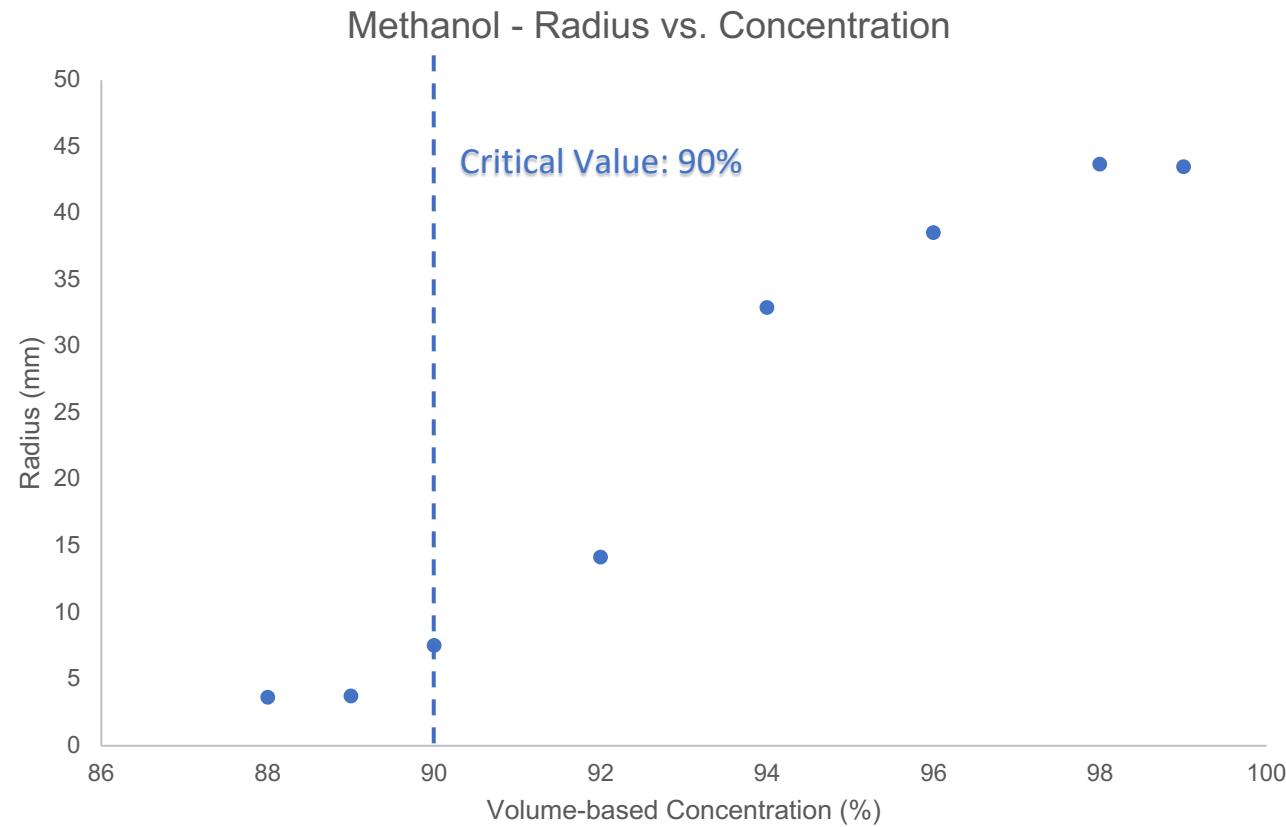


$$\Phi = 0.9$$

Interfacial Tension

Droplet Fragmentation

Critical Value - Vary Concentration

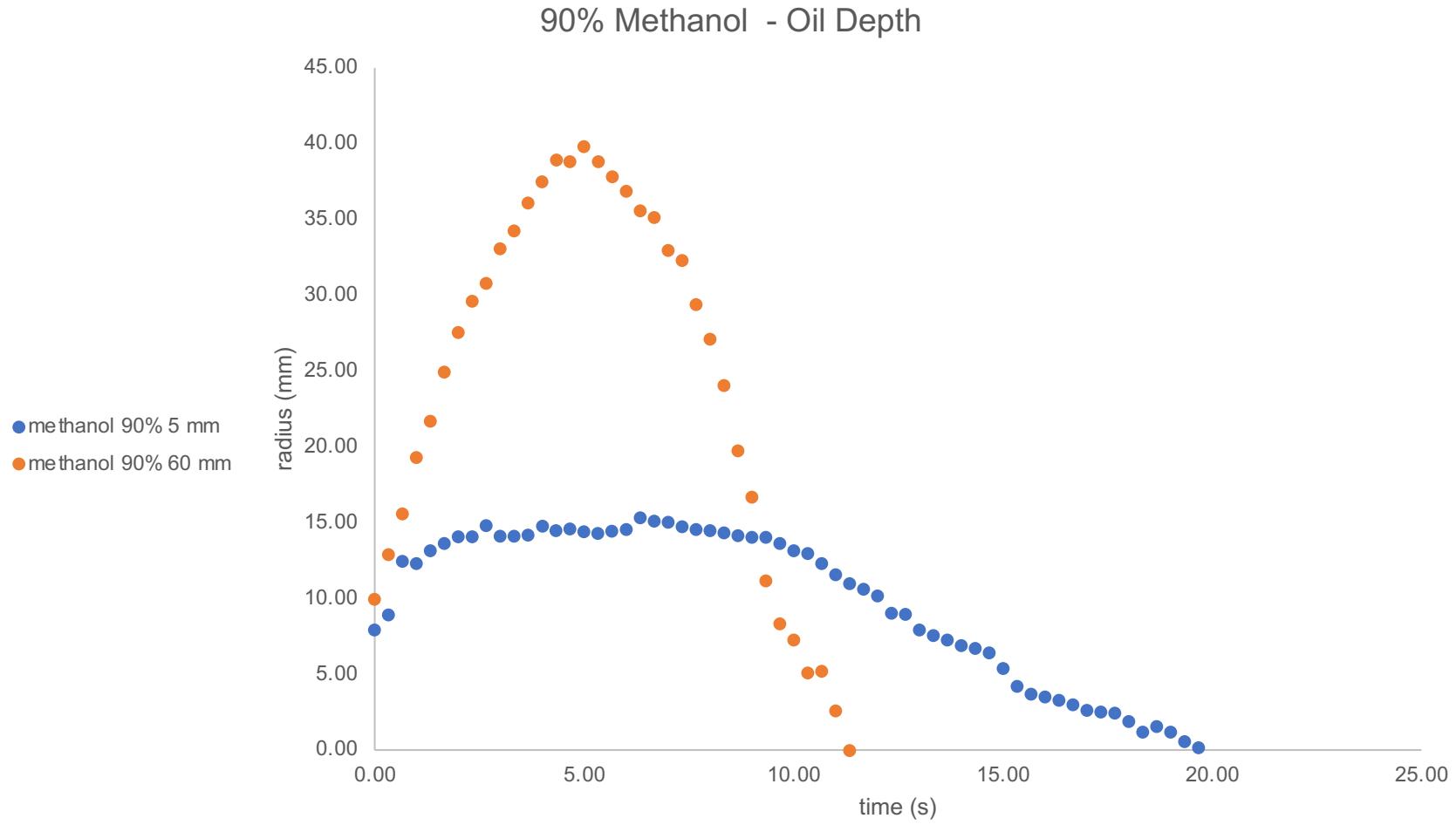


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Vary Oil Depth

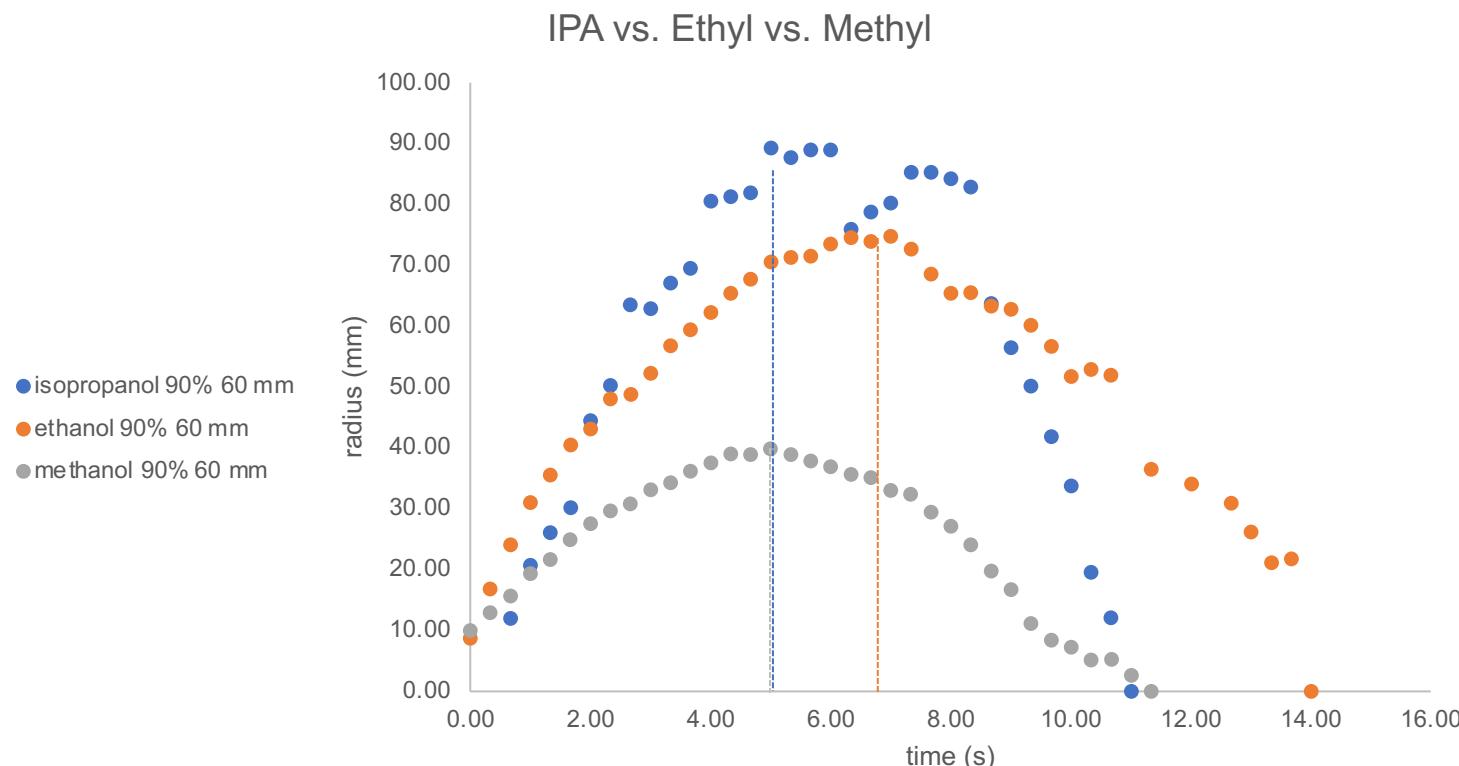


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Vary Solvent

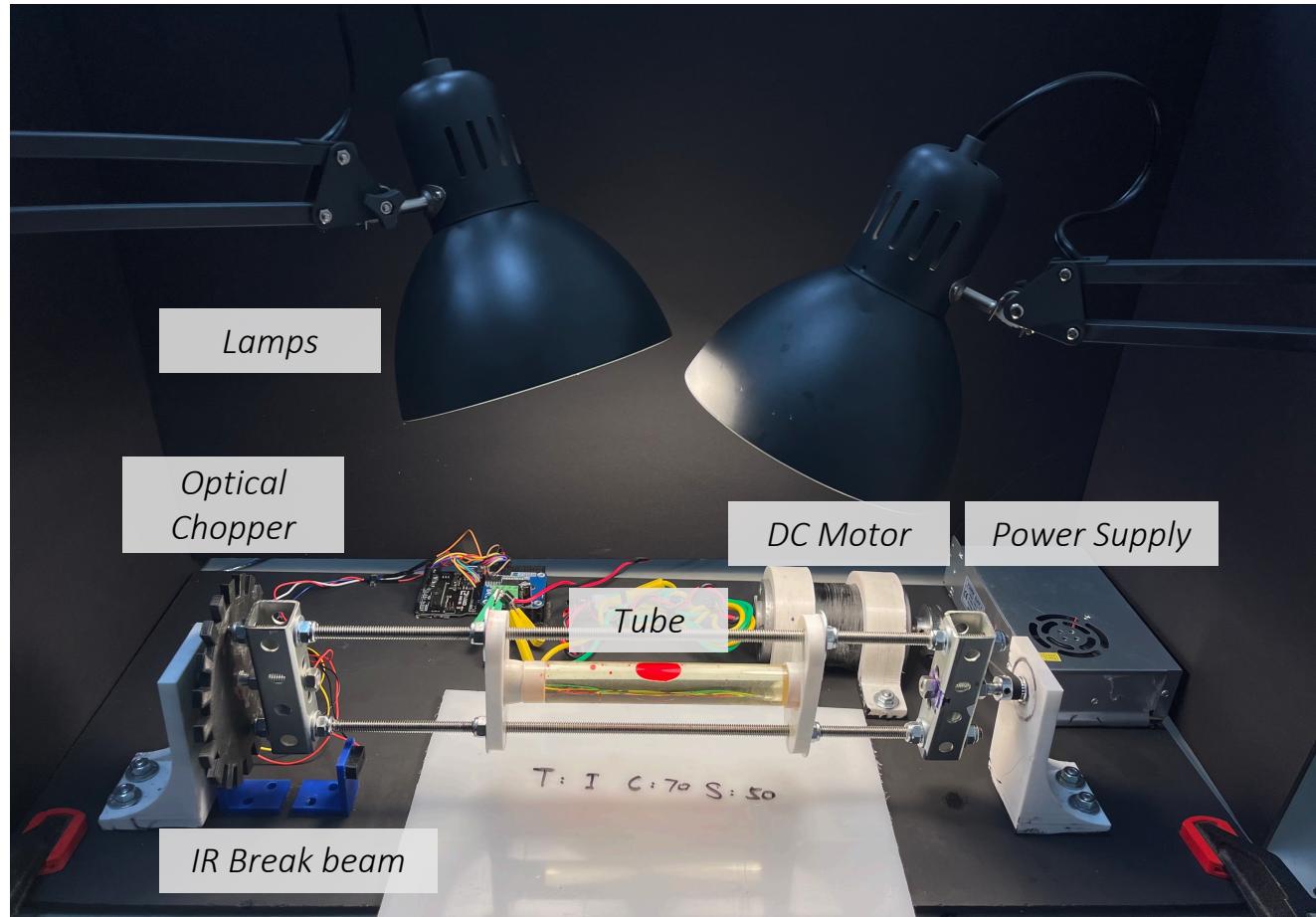


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Interfacial Tension Measurement

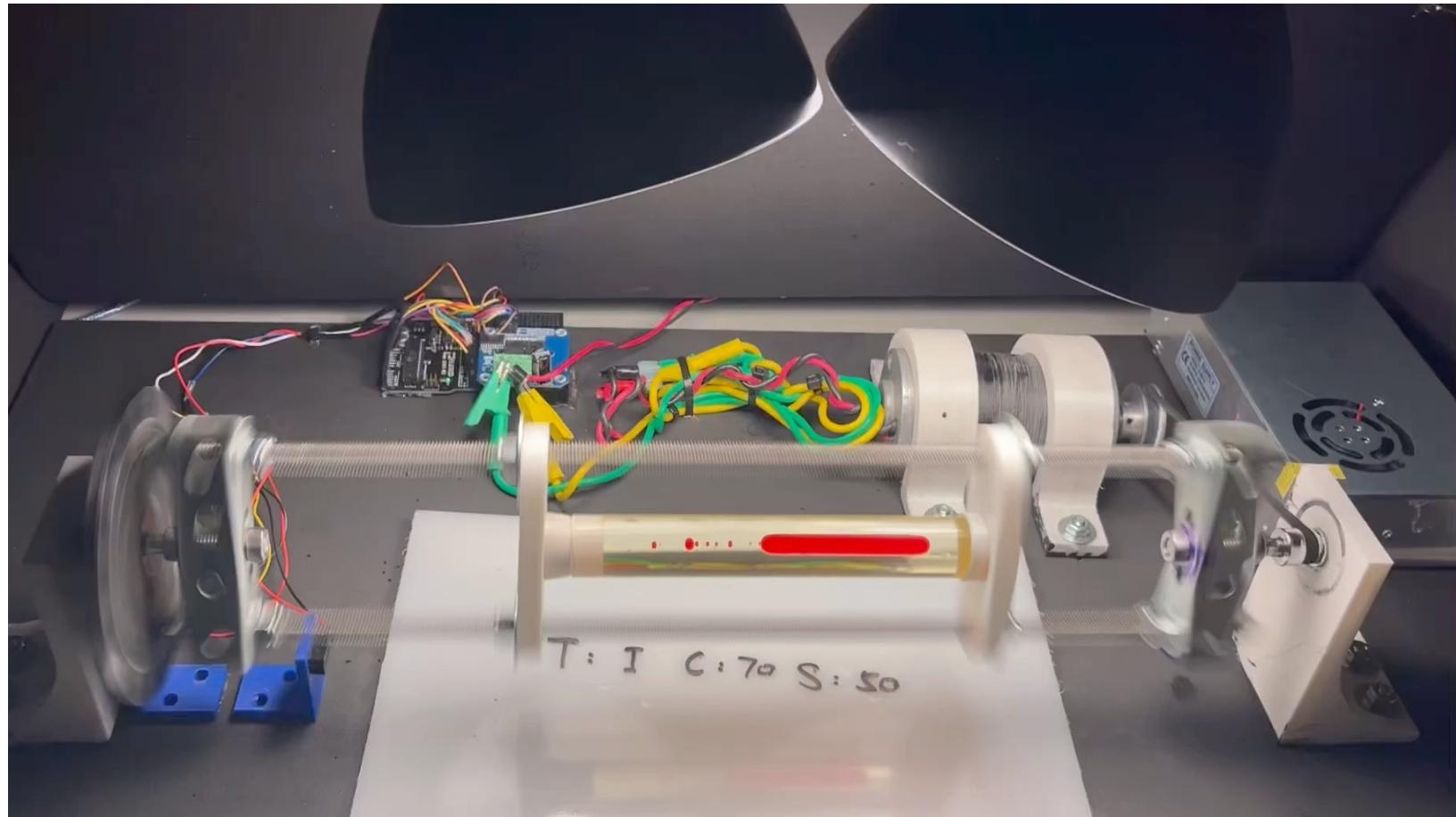


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Interfacial Tension Measurement

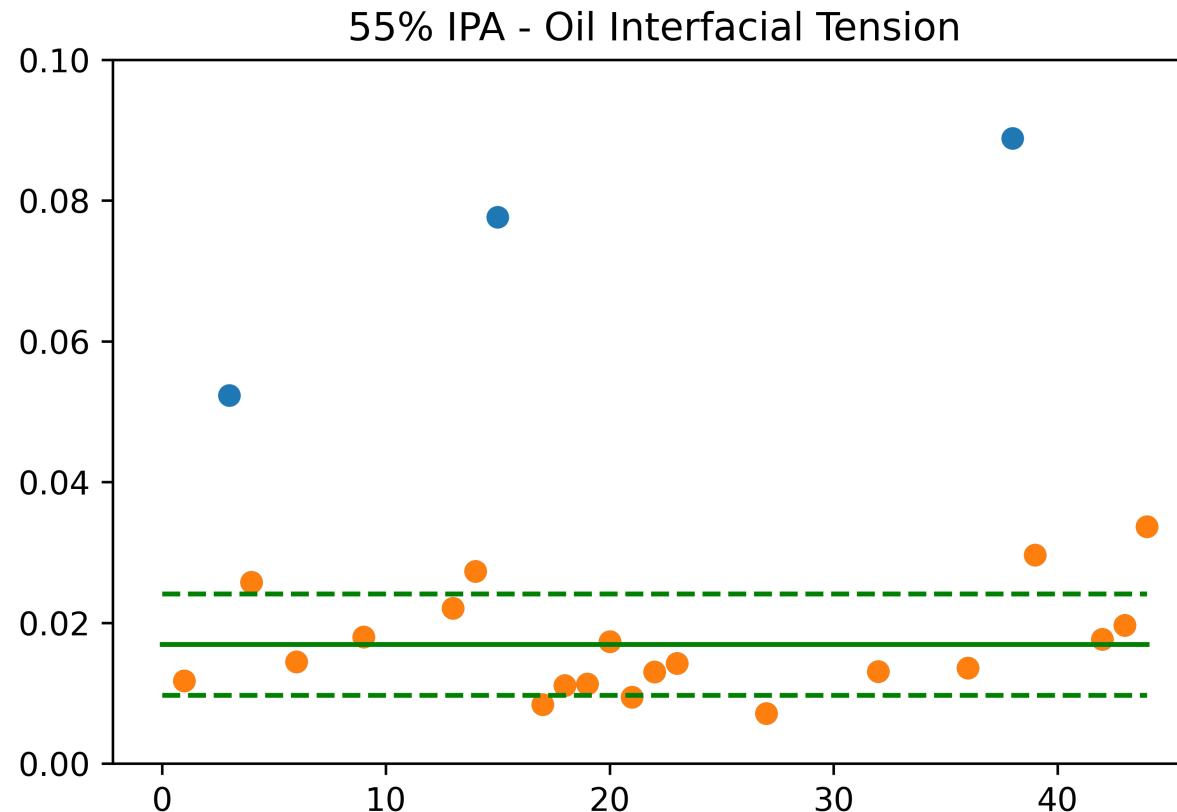


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Interfacial Tension

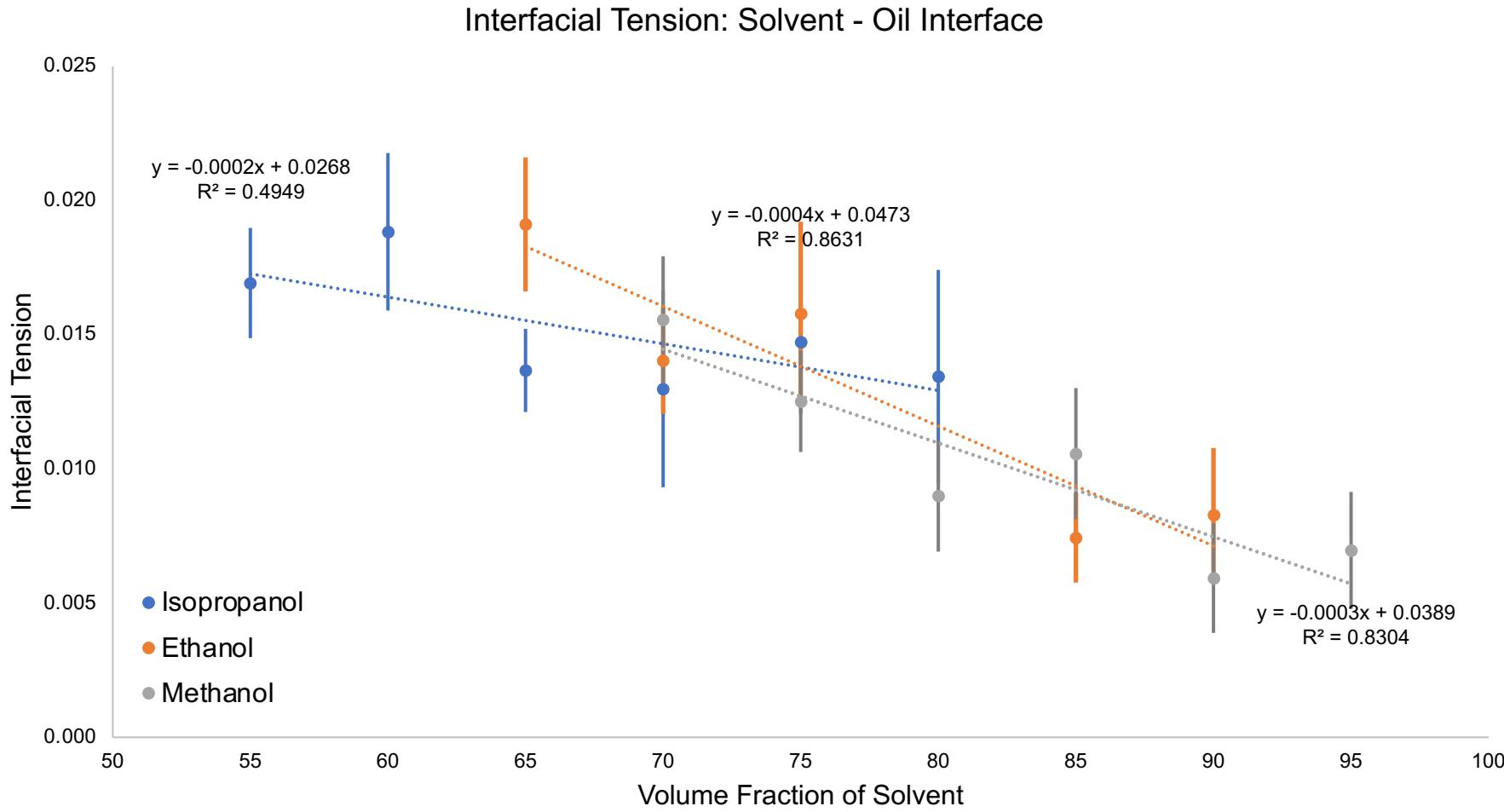


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Interfacial Tension

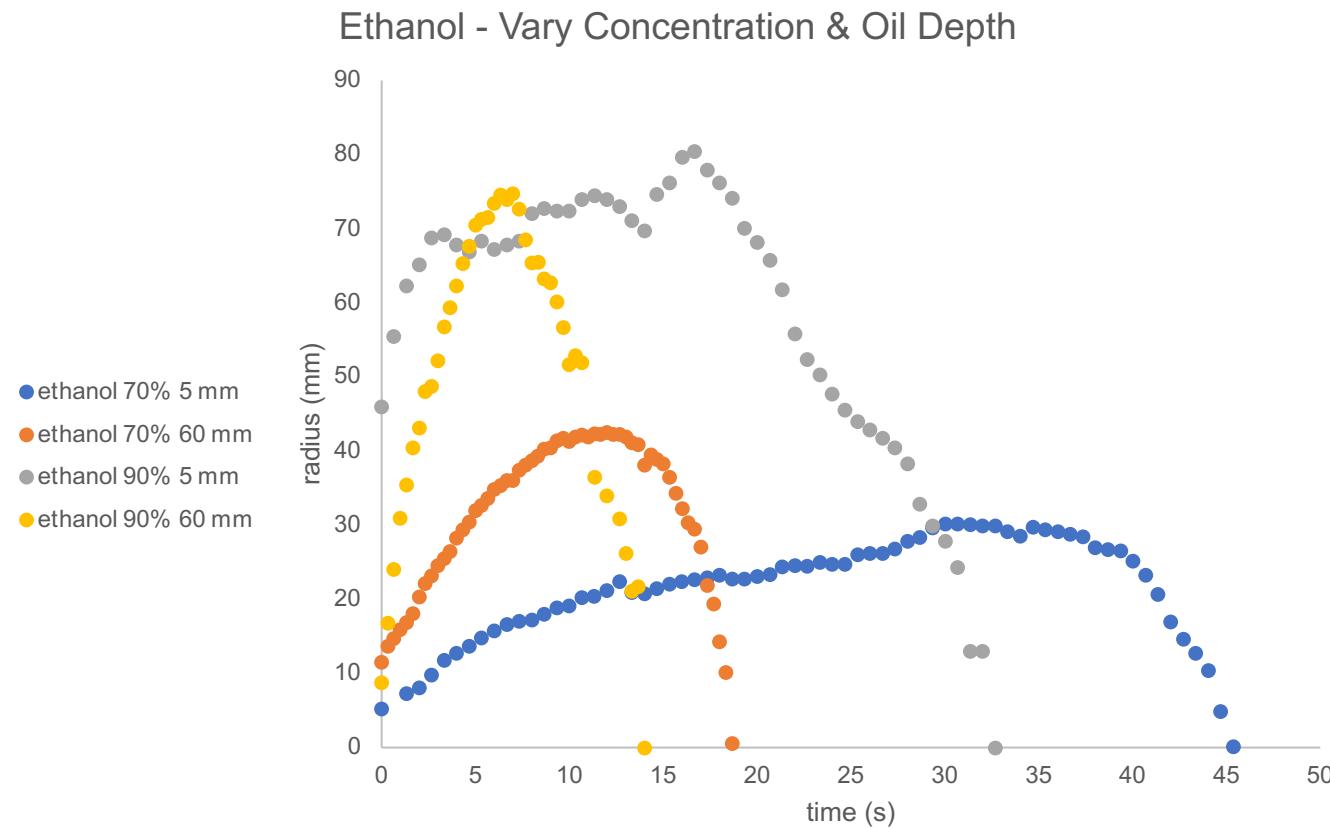


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Experimental Verification

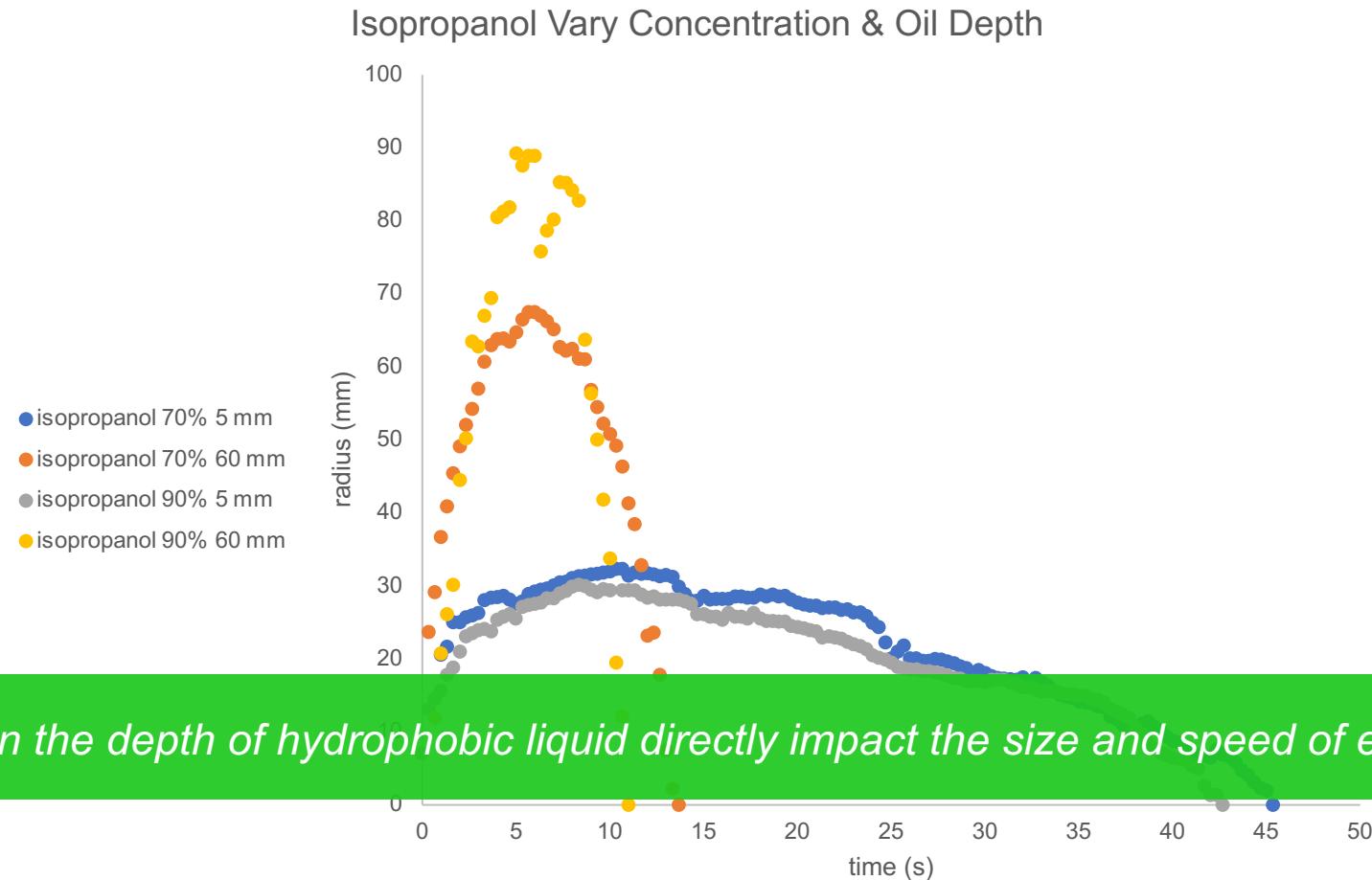


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Experimental Verification



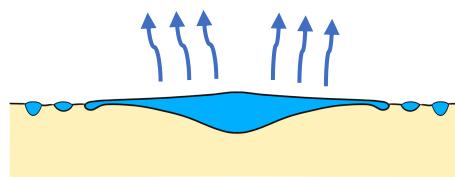
Varies in the depth of hydrophobic liquid directly impact the size and speed of expansion

Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

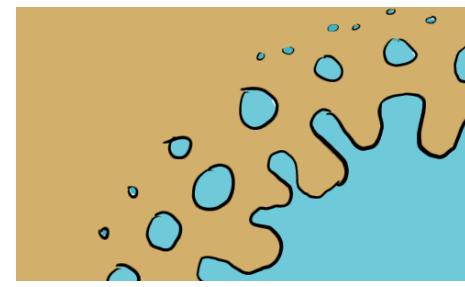
Theoretical Model



Alcohol
Evaporation



Interfacial Tension
& Shear Stress

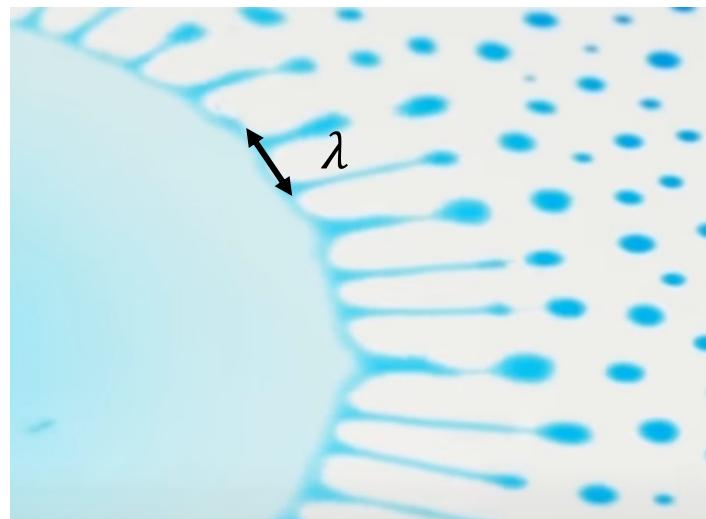


Droplet
Fragmentation

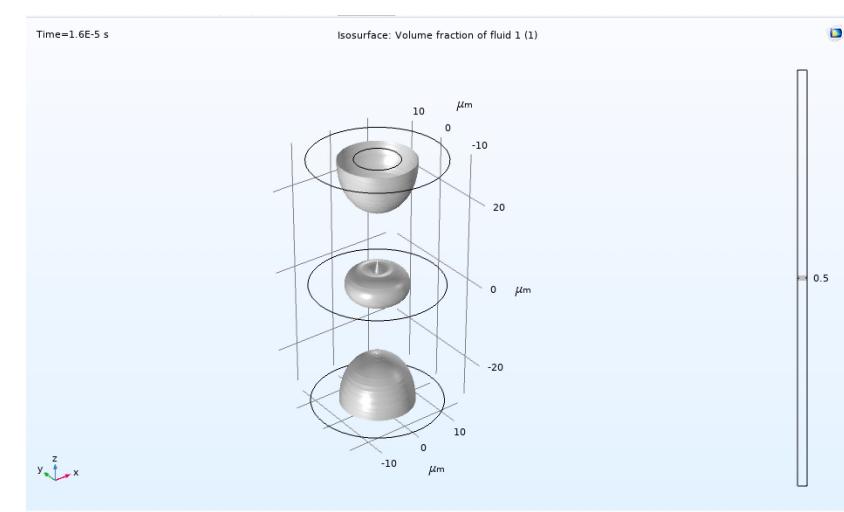
Plateau-Rayleigh Instability

The drop is bounded by a circular rim which destabilizes into small droplets through a process reminiscent of the Plateau-Rayleigh instability of liquid cylinders. Both the wavelength of the instability λ and the size of the resulting droplets strongly depend on the concentration of IPA,

$$\lambda \sim 9.02R_0$$



Alcohol Evaporation



Interfacial Tension

Droplet Fragmentation

Plateau-Rayleigh Instability

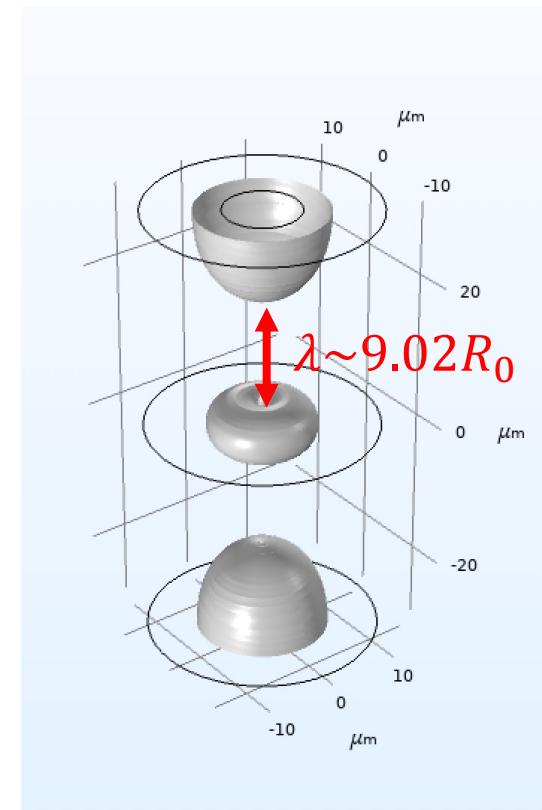
The dependence of the growth rate of the liquid column ω , on wave number k , is:

$$\omega^2 = \frac{\sigma}{\rho R_0^3} k R_0 \frac{I_1 k(R_o)}{I_0(kR_o)} (1 - k^2 R_0^2)$$

For a liquid column (assumed the rim) where σ is the surface tension, ρ is the density, and radius R_o . I_o and I_1 correspond to the solutions of the modified first order Bessel function

And the break up time for the liquid column is,

$$t_{breakup} \sim 2.91 \sqrt{\frac{\rho R_0^3}{\sigma}}$$



Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Scaling of Droplet Size

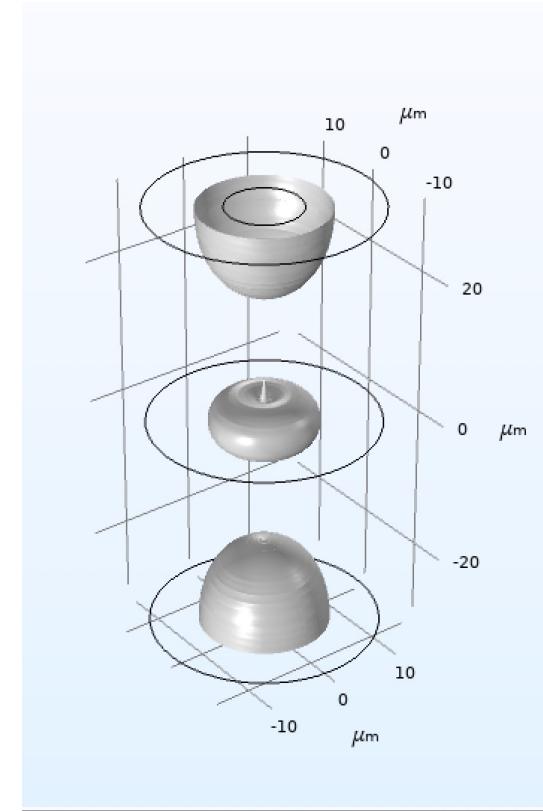
The volume of the main drop,

$$\frac{4}{3}\pi \left[r_{mother\ drop}^3 + \sum r_{daughter\ drop}^3 \right] = \pi r^2 \frac{\omega}{f}$$

While Area is simply, area = πr^2

Substituting the equation derived for the Plateau-Rayleigh instability,

$$V = \pi r^2 \frac{\left(\frac{\sigma}{\rho R_0^3} k R_0 \frac{I_1 k(R_o)}{I_0(kR_0)} (1 - k^2 R_0^2) \right)^{\frac{1}{2}}}{f}$$



As shown above, diameter of ejected droplets, $\sim 20\mu\text{m} - 0.5\text{mm}$

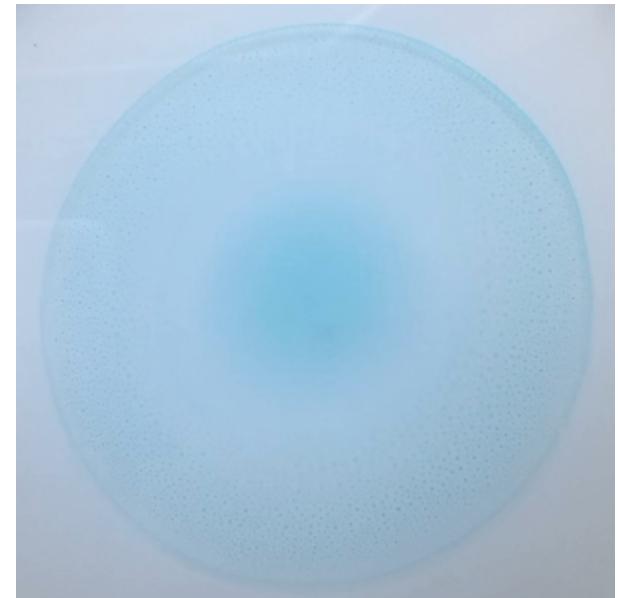
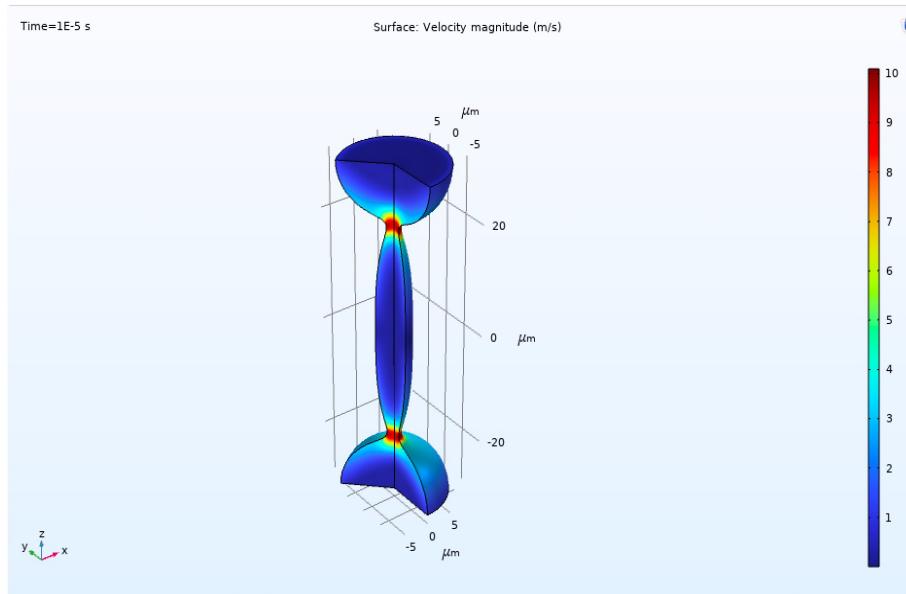
Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Droplet Fragmentation

*Surface tension balances the Marangoni velocity at the rim, and fragmentation occurs,
In which a bulge is created due to this effect*



Alcohol Evaporation

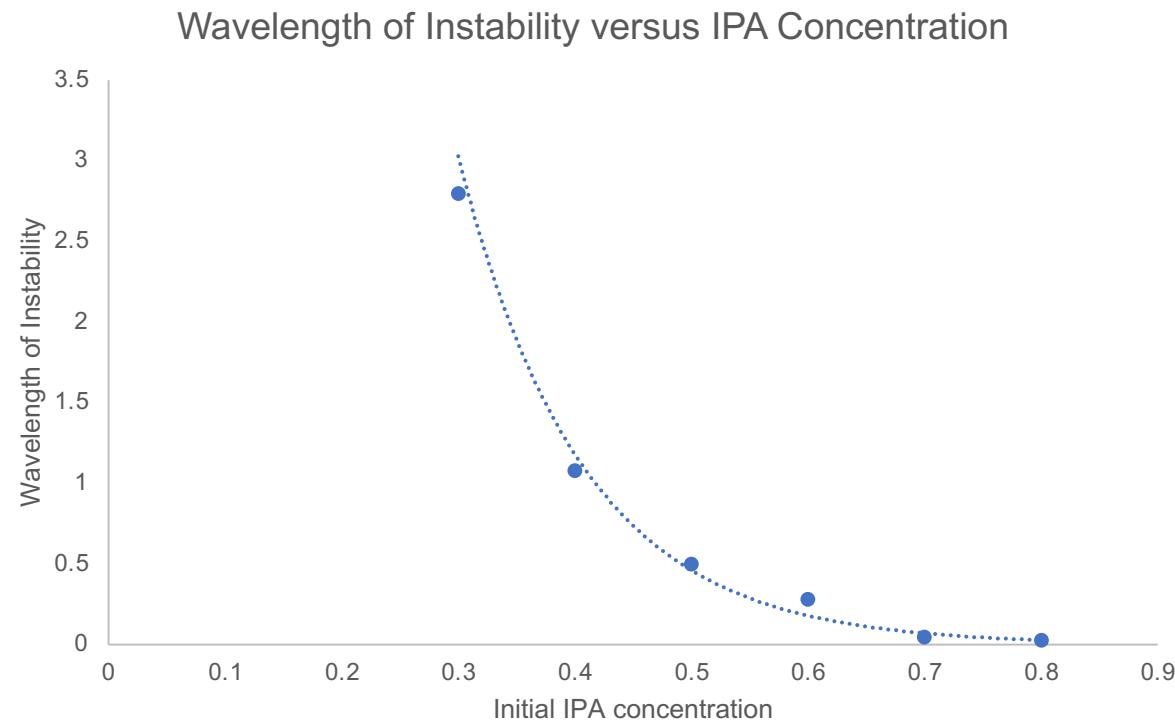
Interfacial Tension

Droplet Fragmentation

Simulation Results

COMSOL microfluidics plateau-Rayleigh instability simulation results

For the wavelength of instability as initial concentration is inputted,



Alcohol Evaporation

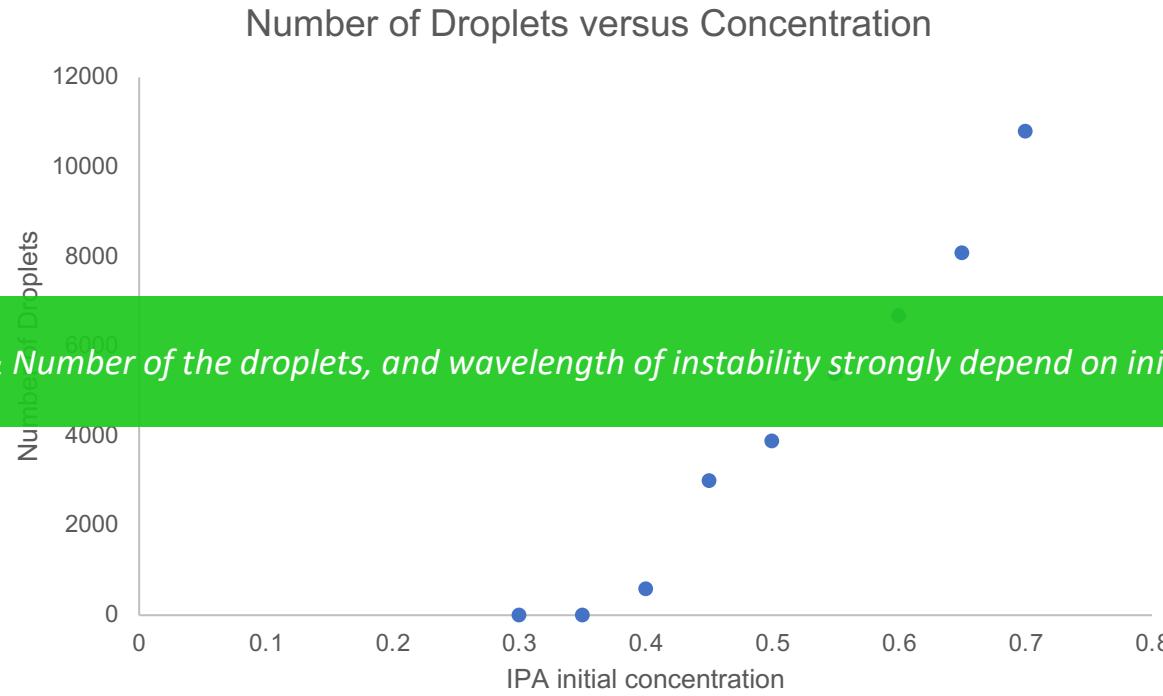
Interfacial Tension

Droplet Fragmentation

Simulation Results

COMSOL microfluidics plateau-Rayleigh instability simulation results

For the number of droplets as initial concentration is inputted,



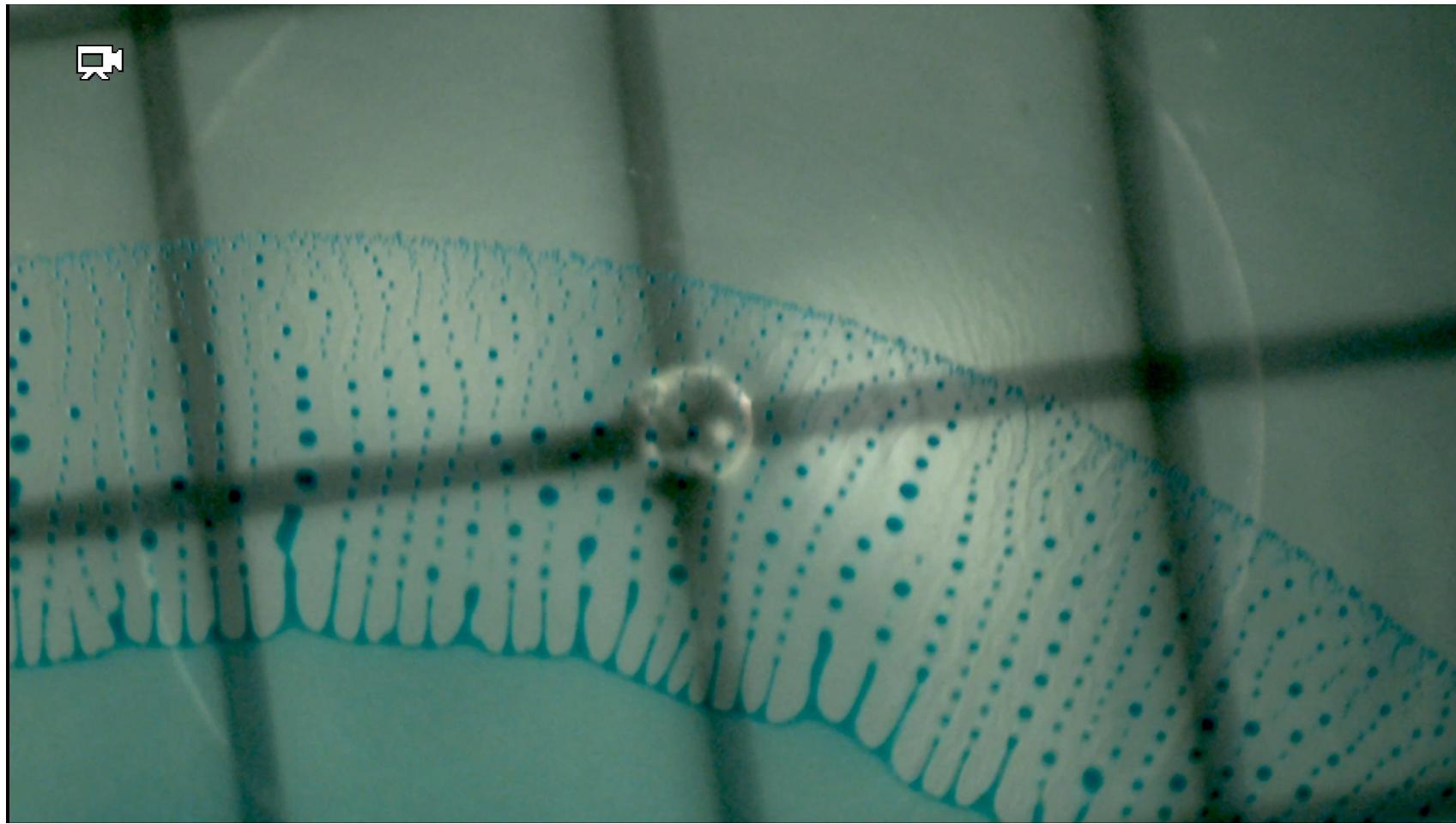
Size & Number of the droplets, and wavelength of instability strongly depend on initial IPA concentrationz

Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Microscopy – 50% Isopropanol

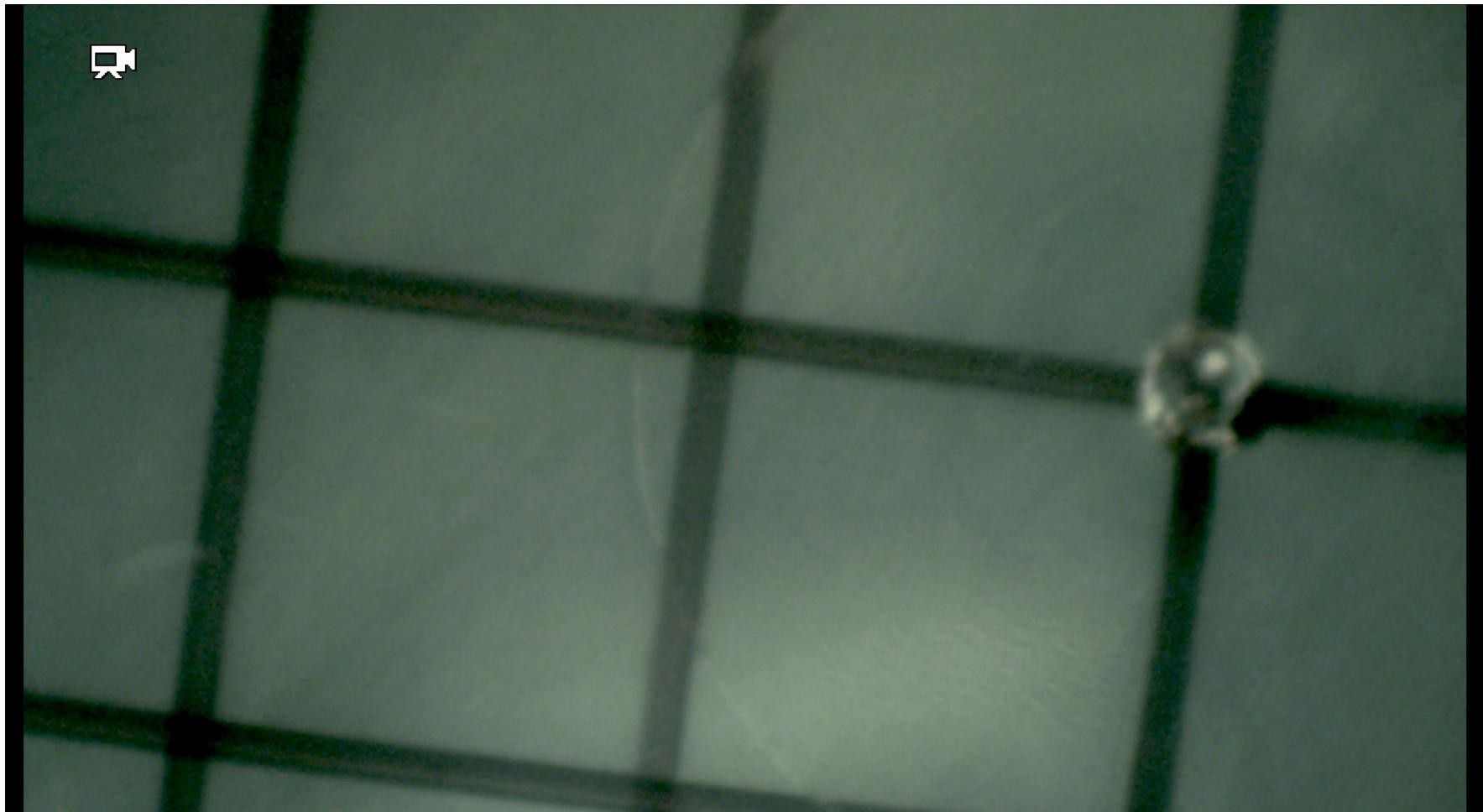


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Microscopy – 70% Isopropanol

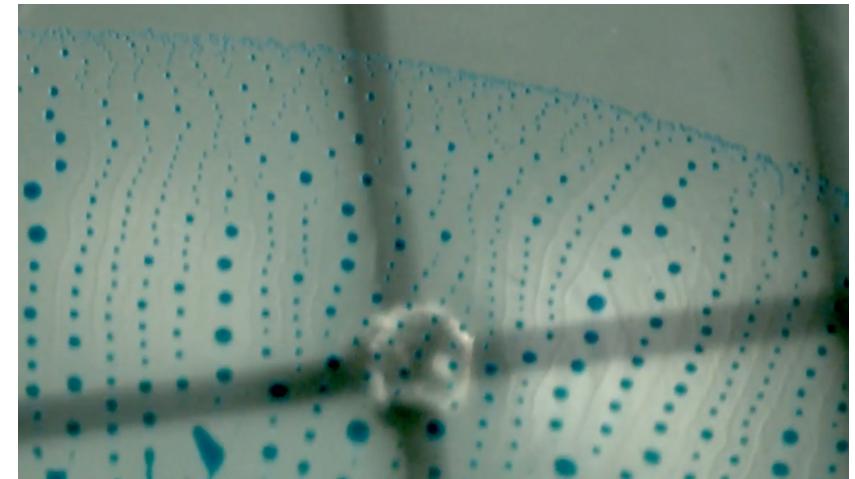
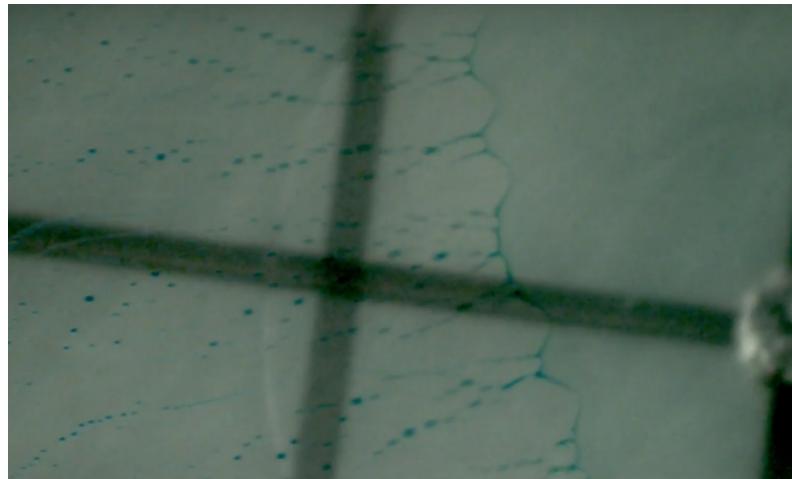


Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Droplet Fragmentations



Higher initial concentration of alcohol in the mother droplet lead to the increase in fragmentation size and number of fragmentation.

70% isopropanol

50% isopropanol

Alcohol Evaporation

Interfacial Tension

Droplet Fragmentation

Key Parameters

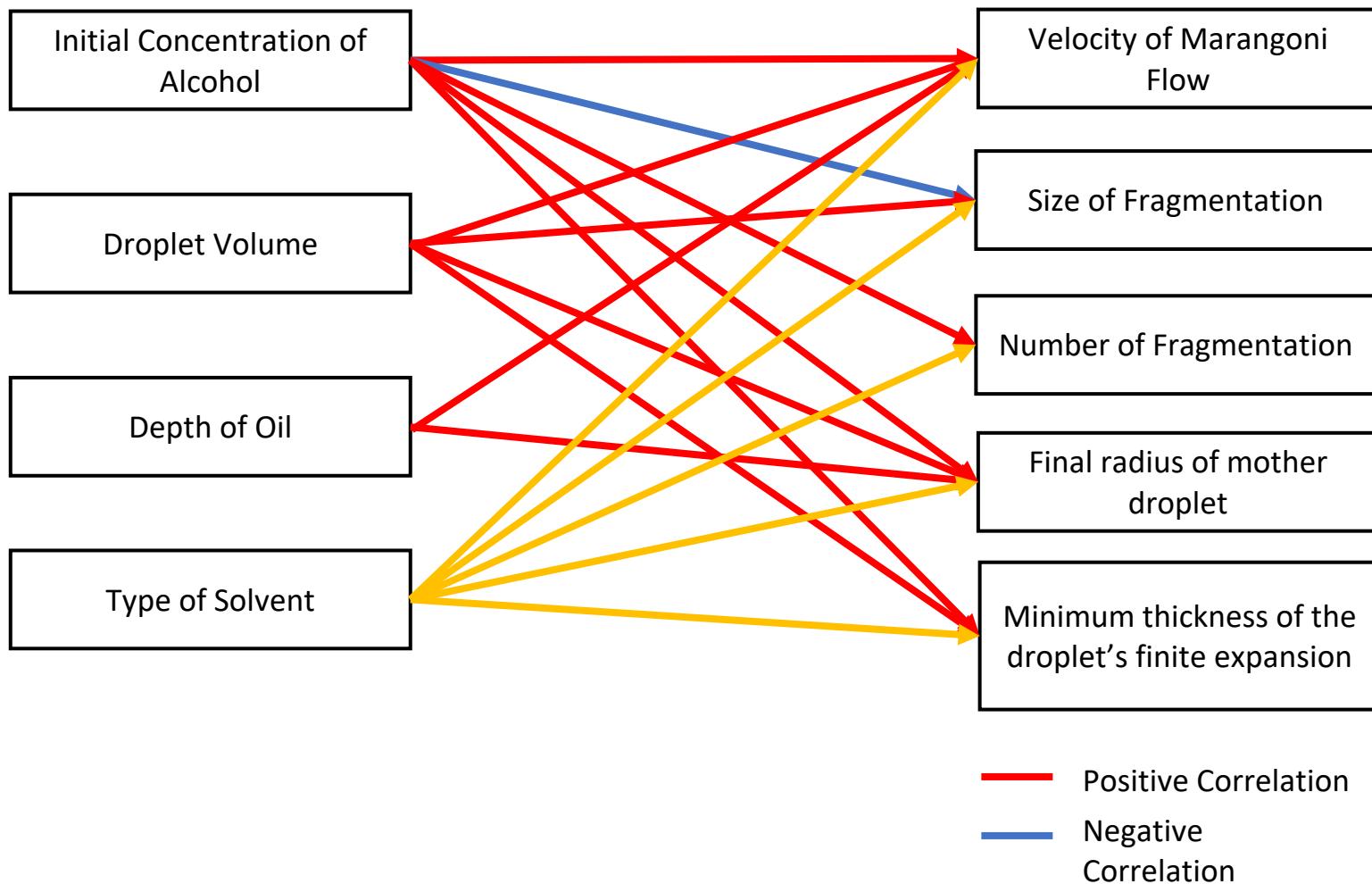
Introduction

Experimental Setup

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Key Parameters

Conclusion

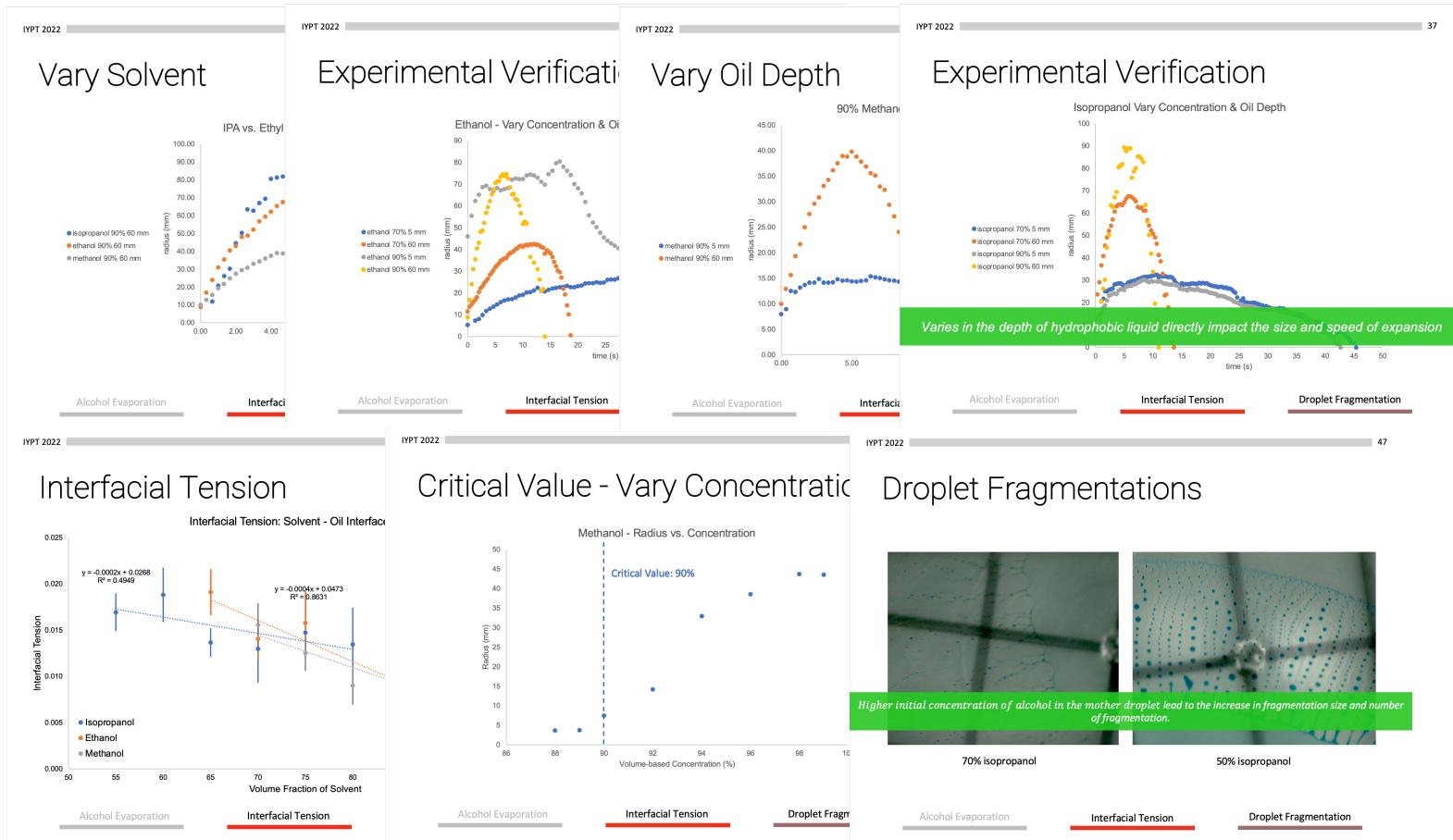


Conclusion

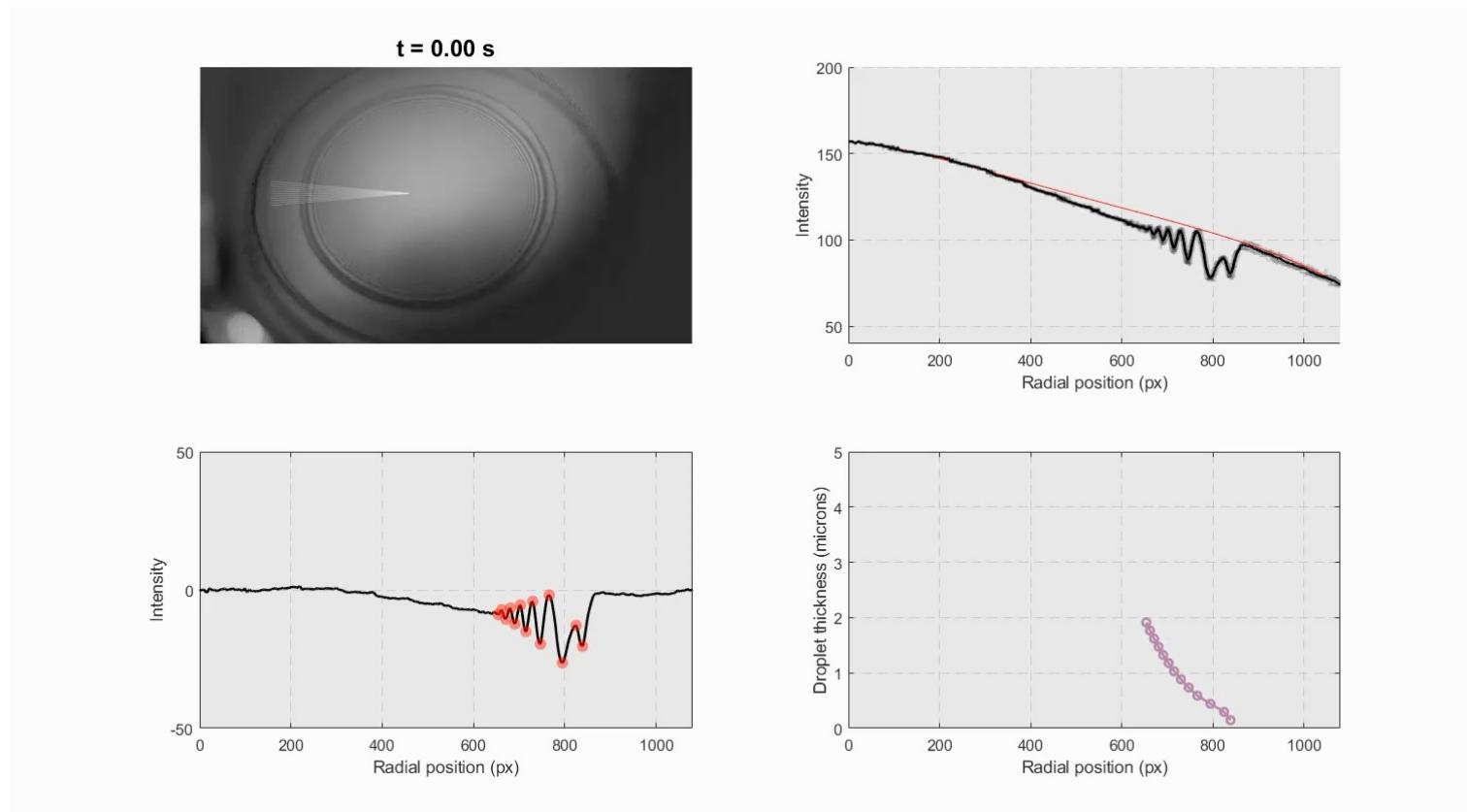
"When a drop of a water mixture (e.g. water-alcohol) is deposited on the surface of a hydrophobic liquid (e.g. vegetable oil), the resulting drop may sometimes fragment into smaller droplets. Investigate the parameters that affect the fragmentation and the size of the final droplets."

Solvent concentration is the **most relevant** parameter to the problem, while other parameters like **hydrophobic liquid surface tension** and **droplet size** have an impact as well. **Fragmentation** occurs through the **surface tension's** forces overcoming **intermolecular bonds**, and **more destabilization** leads to **more and smaller droplets**.

Conclusion



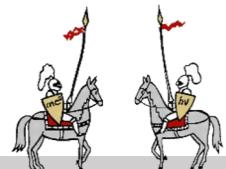
Further Insights



References

- [1] Keiser, L., Bense, H., Colinet, P., Bico, J., Reyssat, E., 2017. Marangoni Bursting: Evaporation-Induced Emulsification of Binary Mixtures on a Liquid Layer. *Physical Review Letters* 118.. doi:10.1103/physrevlett.118.074504
- [2] G. Durey, H. Kwon, Q. Magdelaine, M. Casiulis, J. Mazet, L. Keiser, H. Bense, P. Colinet, J. Bico, and E. Reyssat, *Physical Review Fluids* 3, (2018).
- [3] R. A. Lopez De La Cruz, C. Diddens, X. Zhang, and D. Lohse, *Journal of Fluid Mechanics* 923, (2021).
- [4] H. Machrafi, A. Rednikov, P. Colinet, and P. Dauby, *Journal of Colloid and Interface Science* 349, 331 (2010).
- [5] J. R. Picardo, T. G. Radhakrishna, and S. Pushpavanam, *Journal of Fluid Mechanics* 793, 280 (2016).
- [6] Kimata, M., Terashima, T., Kurita, N., Satsukawa, H., Harada, A., Kodama, K., Sato, A., Imai, M., Kihou, K., Lee, C. H., Kito, H., Eisaki, H., Iyo, A., Saito, T., Fukazawa, H., Kohori, Y., Harima, H., & Uji, S. (2010, December 08). Quasi-two-dimensional Fermi surfaces and coherent interlayer transport in $\{\mathrm{kfe}\}_2\{\mathrm{as}\}_2$. *Physical Review Letters*. Retrieved May 28, 2022, from <https://journals.aps.org/prl/abstract/10.1103/PhysRevLett.105.246403>
- [7] Kimata, M., Terashima, T., Kurita, N., Satsukawa, H., Harada, A., Kodama, K., . . . Uji, S. (2010, December 08). Quasi-two-dimensional Fermi surfaces and coherent interlayer transport in $\{\mathrm{kfe}\}_2\{\mathrm{as}\}_2$. *Physical Review Letters*. Retrieved May 28, 2022, from <https://journals.aps.org/prl/abstract/10.1103/PhysRevLett.105.246403>
- [8] A. D'Aubeterre, R. Da Silva, and M. E. Aguilera, *International Communications in Heat and Mass Transfer* 32, 677 (2005).

Thank you for listening



Appendix

Appendix

Interfacial Tension Calculation

```
1 import numpy as np
2 import matplotlib.pyplot as plt
3
4 import glob
5 print("Using glob.glob()")
6 files = glob.glob("//Users/dongyuezhen/Desktop/data_cal_copy/55 IPA/*.csv",
7                 recursive = True)
8
9 speed = np.zeros(0)
10 for file in files:
11     #print(float(str.split(file, '_')[4]))
12     speed = np.append(speed, [float(str.split(file, '_')[4])])
13     #print(file)
14
15
16 d0 = 0.92
17 dm = 0.91340
18 dd = d0-dm
19 omega = speed
20
21 #initialize H,G arrays
22 H = np.zeros(len(speed))
23 G = np.zeros(len(speed))
24 x = np.zeros(len(speed))
25
26 for i in range(len(files)):
27     x, y = np.loadtxt(files[i], delimiter = ',', unpack = True)
28     x = x/1000
29     x0 = (np.max(x) - np.min(x))/2
30     x[i] = x0
31     H[i] = (dd*omega[i]**2)/4
32     G[i] = (3*x0/2)*((dd*omega[i]**2)/4)**(1/3)
33     #plt.show()
34     #print(x , speed)
35 print('Data File Number = '+str(len(speed)+1))
36
37
38 #initialize array to compare every combination of two points
39 combination = len(speed)*(len(speed)-1)/2
40 index = np.zeros((int(combination), 2))
```

Interfacial Tension Calculation

```
43 counter = 0
44 for i in range(len(speed)):
45     for j in range(len(speed)):
46         if(i!=j and i<j):
47             index[counter] = [i,j]
48             counter += 1
49
50 gamma = np.zeros(int(combination))
51
52 for i in range(int(combination)):
53     i1 = int(index[i][0])
54     i2 = int(index[i][1])
55     gamma[i] = G[i1] - H[i1]*((G[i1]-G[i2])/(H[i1]-H[i2]))
56
57
58 #gamma = filter_until(gamma,0.1)
59
60 m = np.arange(0,len(gamma), 1)
61 print(gamma)
62
63 gamma_clip = gamma
64 m_clip = m
65
66
67 m_clip = m_clip[gamma_clip > 0]
68 gamma_clip = gamma_clip[gamma_clip > 0]
69 m_clip = m_clip[gamma_clip <= 4*10**(-2)]
70
71
72 gamma_clip = gamma_clip[gamma_clip <= 4*10**(-2)]
73
74
75
76 print(gamma_clip)
77
78 gammabar = np.average(gamma_clip)
79 gammasd = np.std(gamma_clip)
80 print(gammabar, gammasd)
81 plt.plot(m,gamma, 'o')
82 plt.plot(m_clip, gamma_clip, 'o')
83
84 plt.plot([m.min(), m.max()], [gammabar, gammabar], 'g')
```

Interfacial Tension Calculation

```
55 gamma_12 = G[11] - m[11] - ((G[11]-G[12])/(m[11]-m[12]))
56
57
58 #gamma = filter_until(gamma,0.1)
59
60 m = np.arange(0,len(gamma), 1)
61 print(gamma)
62
63 gamma_clip = gamma
64 m_clip = m
65
66
67 m_clip = m_clip[gamma_clip > 0]
68 gamma_clip = gamma_clip[gamma_clip > 0]
69 m_clip = m_clip[gamma_clip <= 4*10**(-2)]
70
71
72 gamma_clip = gamma_clip[gamma_clip <= 4*10**(-2)]
73
74
75
76 print(gamma_clip)
77
78 gammabar = np.average(gamma_clip)
79 gammasd = np.std(gamma_clip)
80 print(gammabar, gammasd)
81 plt.plot(m,gamma, 'o')
82 plt.plot(m_clip, gamma_clip, 'o')
83
84 plt.plot([m.min(), m.max()], [gammabar, gammabar], 'g')
85 plt.plot([m.min(), m.max()], [gammabar+gammasd, gammabar+gammasd], 'g--')
86 plt.plot([m.min(), m.max()], [gammabar-gammasd, gammabar-gammasd], 'g--')
87 plt.ylim(0,1e-1)
88 plt.title('55% IPA - Oil Interfacial Tension')
89 plt.savefig('55%_IPA_interfacial', facecolor = "white", dpi = 600)
90 plt.show()
91
92
93
94
95
96
97
```