

MASSACHUSETTS INSTITUTE OF TECHNOLOGY  
Department of Chemistry

5.311 DATA SUMMARY SHEET FALL 2005

Student name: \_\_\_\_\_

TA name: \_\_\_\_\_

Group: A / B (Please circle)

Section: MW / TR (Please circle)

**EXPERIMENT 1: KINETICS**

molar absorptivity ( $\epsilon$ )	
$k_{\text{obs}}$	
$k_o$	
$k_{\text{real}}$	
$Z_1Z_2$	
Correlation coefficient (R) Debye-Huckel plot	
temperature (t)	
activation energy ( $E_a$ )	
Arrhenius parameter (A)	
Correlation coefficient (R) Arrhenius plot	

MASSACHUSETTS INSTITUTE OF TECHNOLOGY  
Department of Chemistry

5.311 DATA SUMMARY SHEET FALL 2005

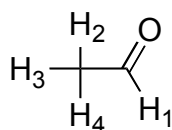
Student name: \_\_\_\_\_

TA name: \_\_\_\_\_

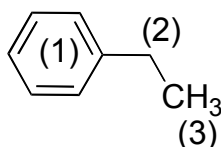
Group: A / B (Please circle)

Section: MW / TR (Please circle)

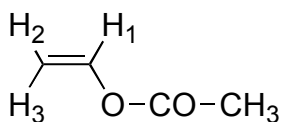
**EXPERIMENT 2: Nuclear Magnetic Resonance**



Proton	Chemical Shift (ppm)	Multiplicity	Integration	J-Couplings
H <sub>1</sub>				
H <sub>2</sub>				
H <sub>3</sub>				
H <sub>4</sub>				



Proton	Chemical Shift (ppm)	Multiplicity	Integration	J-Couplings
H <sub>1</sub>				
H <sub>2</sub>				
H <sub>3</sub>				



Proton	Chemical Shift (ppm)	Multiplicity	Integration	J-Couplings
H <sub>1</sub>				
H <sub>2</sub>				
H <sub>3</sub>				





MASSACHUSETTS INSTITUTE OF TECHNOLOGY  
Department of Chemistry

5.311 DATA SUMMARY SHEET FALL 2005

Student name: \_\_\_\_\_

TA name: \_\_\_\_\_

Group: A / B (Please circle)

Section: MW / TR (Please circle)

**EXPERIMENT 3: Unknown Aromatic Carboxylic Acid**

**A.**

Unknown number	Experimental melting point of the acid	Equivalent weight of the acid	pKa of the acid	Identity of the carboxylic acid
	Crude:			
	Recrystallized:			

Yield of acid	Theoretical yield	Percent yield
Crude:		
Recrystallized:	*	*

\*base on the mass of crude used in the recrystallization. Mass of crude used \_\_\_\_\_

**B.**

	Titration 1	Titration 2	Titration 3
Calculated molarity of the sodium hydroxide solution for each titration with error propagation			

Report any additional titrations on an additional page.

	reading the buret	mass	endpoint location
Estimated uncertainty in measurement			

95% Confidence limits: \_\_\_\_\_

Average propagated error: \_\_\_\_\_

Concentration of NaOH: \_\_\_\_\_



C.

	Titration 1	Titration 2	Titration 3
Calculated equivalent weight from each titration with error propagation			

Report any additional titrations on an additional page.

95% Confidence limits of mean equivalent weight: \_\_\_\_\_

Average propagated error in equivalent weight: \_\_\_\_\_

D.

	Titration 1	Titration 2	Titration 3
Calculated $V_e$ from each titration with error propagation			

	Titration 1	Titration 2	Titration 3
Calculated pKa from each titration			

Report any additional titrations on an additional page.

pKa: \_\_\_\_\_

95% Confidence limits of pKa: \_\_\_\_\_



MASSACHUSETTS INSTITUTE OF TECHNOLOGY  
Department of Chemistry

5.311 DATA SUMMARY SHEET FALL 2005

Student name: \_\_\_\_\_

TA name: \_\_\_\_\_

Group: A / B (Please circle)

Section: MW / TR (Please circle)

**EXPERIMENT 4: Ferrocene and Acetylferrocenes**

Yield of ferrocene from synthesis (crude)	Limiting reagent	Theoretical yield	Percent yield

	Mass of pure ferrocene	Theoretical yield (based on the mass used in technique)	Percent yield (based on the mass used in technique)
Sublimation			
Recrystallization			

mass [emim]I	mass AlCl <sub>3</sub>	mass ferrocene	volume acetic anhydride	time stirring	crude yield

Observations TLC (including R<sub>f</sub> values, solvent systems): Please provide in a table on the reverse.

GCMS results	Ferrocene (FW=186)	Acetylferrocene (FW=228)	Diacetyl ferrocene (FW=270)

Solvent system used for column chromatography: \_\_\_\_\_

	Ferrocene (unreacted)	Acetylferrocene	Diacetyl ferrocene	Diacetyl ferrocene
Yield				
Percent Yield				
Melting Point Range				