

COMPASS_TEMPEST_SGW_2025: Well Test

Stephanie J. Wilson

2025-02-11

Contents

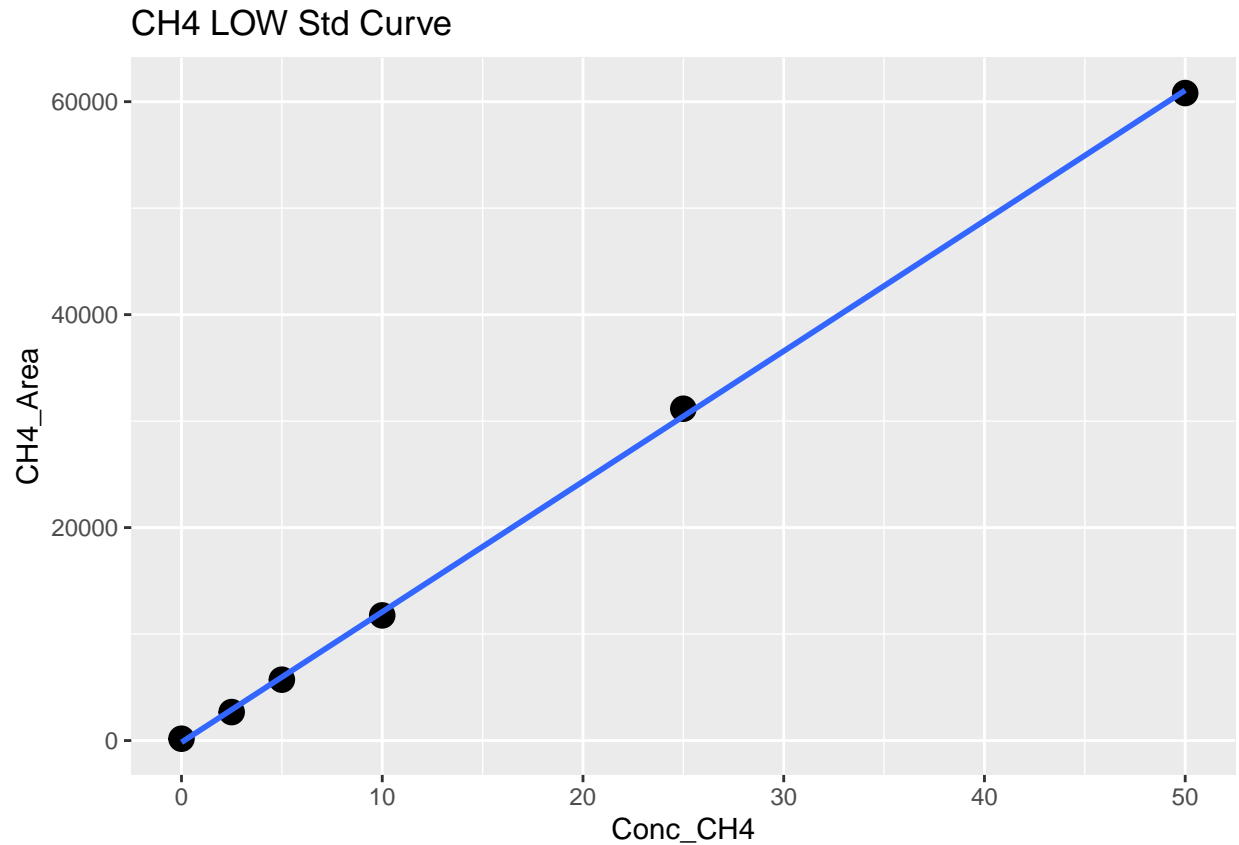
0.1	Read in first data file and assess standard curves	1
0.2	Now calculate the CH4 & CO2 concentrations in ppm	4
0.3	Check the Check Standards	5
0.4	Dilution correct samples	5
0.5	If samples are water calculate gas in water	7
0.6	Write out processed data & slopes	7

##Set Up

0.1 Read in first data file and assess standard curves

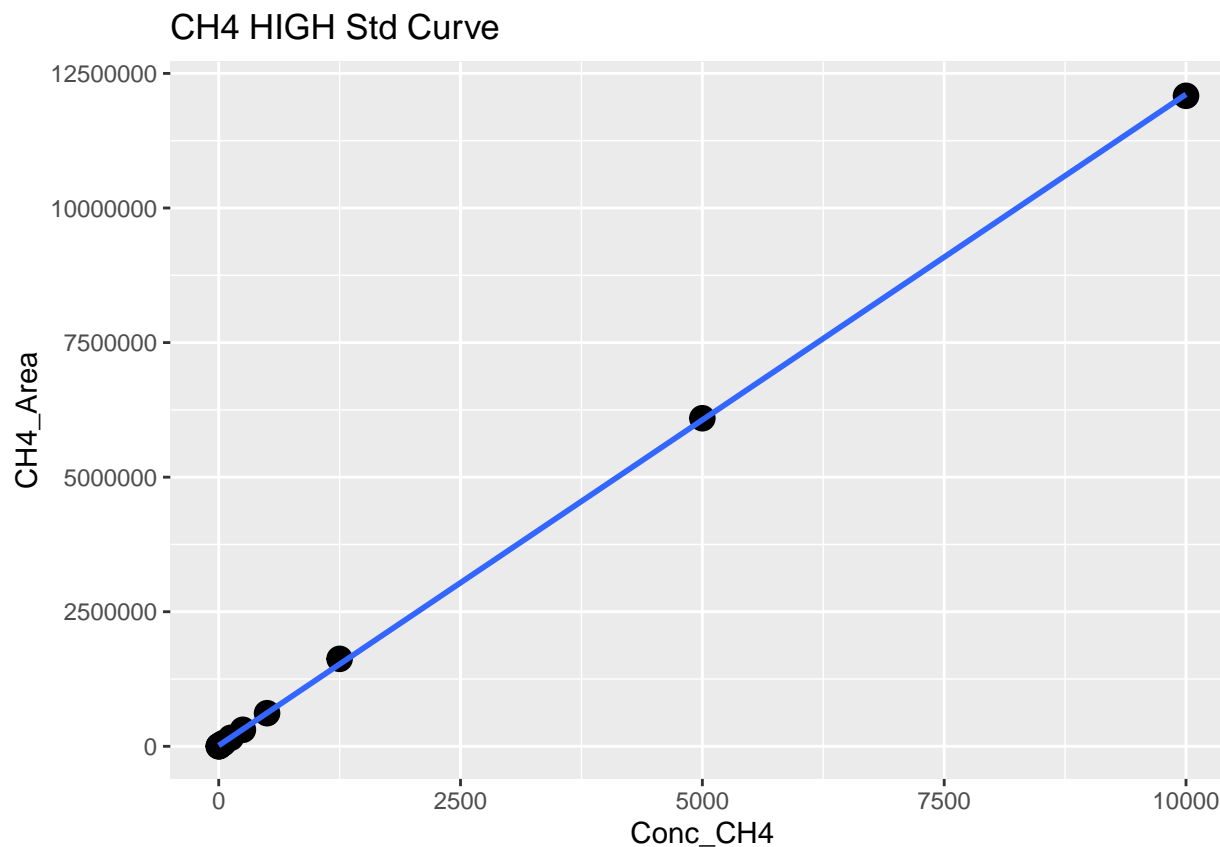
```
##      Machine      User Run_Date Sample_Year Sample_Month Sample_Day
## 1 Shimadzu GC B. Blakley 20250922      NA      <NA>      NA
## 2 Shimadzu GC B. Blakley 20250922      NA      <NA>      NA
## 3 Shimadzu GC B. Blakley 20250922      NA      <NA>      NA
## 4 Shimadzu GC B. Blakley 20250922      NA      <NA>      NA
## 5 Shimadzu GC B. Blakley 20250922      NA      <NA>      NA
## 6 Shimadzu GC B. Blakley 20250922      NA      <NA>      NA
##  Event_Stamp Sample_Time Time_Zone  G_W Sample_ID Sample_Type Conc_CO2
## 1      <NA>      NA      <NA> <NA>   LabAir      Lab air      NA
## 2      <NA>      NA      <NA> <NA>   Blank1      Blank      NA
## 3      <NA>      NA      <NA> <NA>   Blank2      Blank      NA
## 4      <NA>      NA      <NA> <NA>     Oppm      Standard  NA
## 5      <NA>      NA      <NA> <NA>   2.5ppm      Standard  NA
## 6      <NA>      NA      <NA> <NA>     5ppm      Standard  NA
##  Conc_CH4 Conc_N2O CO2_Area CH4_Area Dilution_Factor Hours
## 1      NA      NA      NA      1277      NA      NA
## 2      NA      NA      NA      164      NA      NA
## 3      NA      NA      NA      166      NA      NA
## 4      0.0      NA      NA      172      NA      NA
## 5      2.5      NA      NA      2672     NA      NA
## 6      5.0      NA      NA      5716     NA      NA
```

```
## 'geom_smooth()' using formula = 'y ~ x'
```



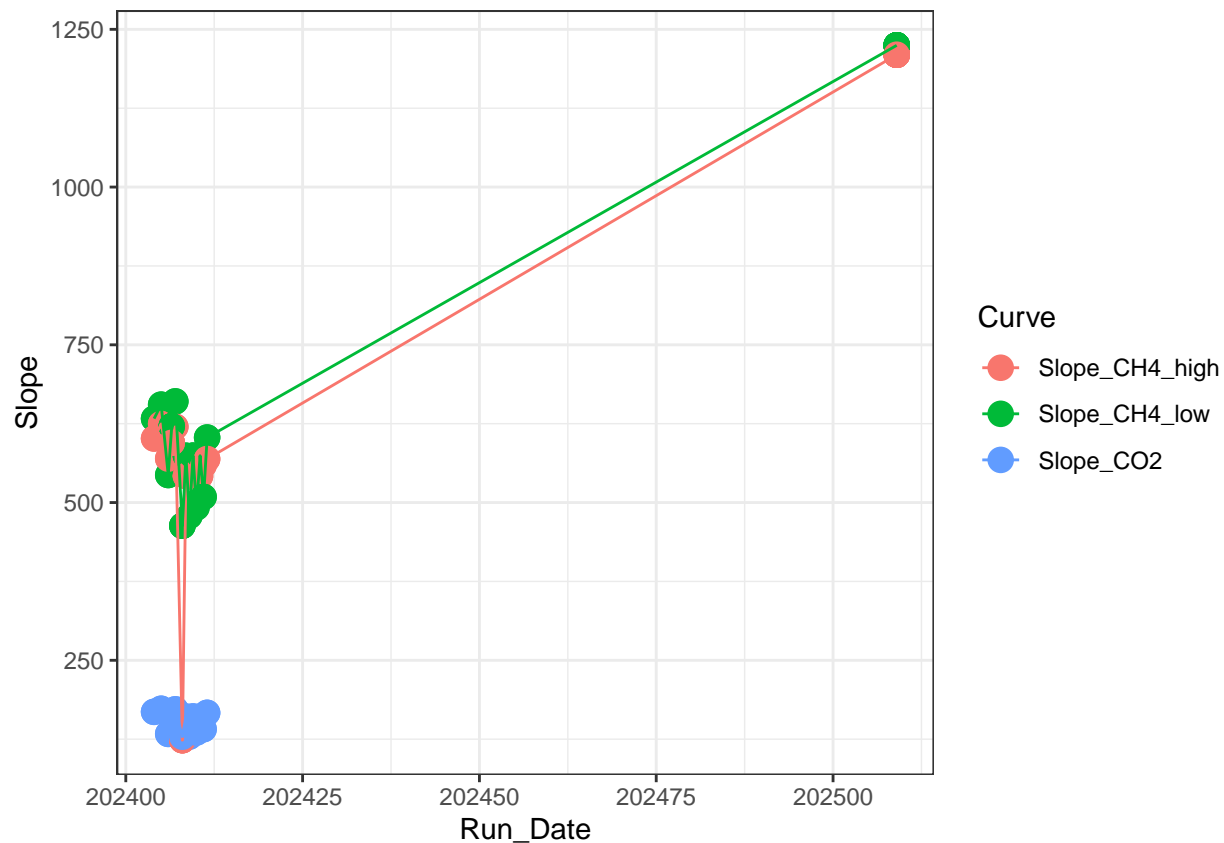
```
##
## Call:
## lm(formula = stds_ch4_low$CH4_Area ~ stds_ch4_low$Conc_CH4)
##
## Residuals:
##      1      2      3      4      5      6
## 337.1 -224.4 -241.9 -327.9  712.0 -255.0
##
## Coefficients:
##              Estimate Std. Error t value Pr(>|t|)
## (Intercept)    -165.1     258.6  -0.639   0.558
## stds_ch4_low$Conc_CH4  1224.6       11.1 110.319 4.05e-08 ***
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 474.9 on 4 degrees of freedom
## Multiple R-squared:  0.9997, Adjusted R-squared:  0.9996
## F-statistic: 1.217e+04 on 1 and 4 DF, p-value: 4.049e-08

## 'geom_smooth()' using formula = 'y ~ x'
```



```
##
## Call:
## lm(formula = stds_ch4$CH4_Area ~ stds_ch4$Conc_CH4)
##
## Residuals:
##      Min       1Q   Median       3Q      Max
## -27445 -14156 -13201  -6134   98453
##
## Coefficients:
##              Estimate Std. Error t value Pr(>|t|)
## (Intercept)    13820.463   11476.617     1.204   0.256
## stds_ch4$Conc_CH4 1209.769     3.529 342.782 <2e-16 ***
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 35680 on 10 degrees of freedom
## Multiple R-squared:  0.9999, Adjusted R-squared:  0.9999
## F-statistic: 1.175e+05 on 1 and 10 DF,  p-value: < 2.2e-16
```

	X	Curve	R2	Slope	Intercept	Run_Date
## 1	1	Slope_CH4_low	0.9996407	633.0314	-320.6965	202404
## 2	2	Slope_CH4_high	0.9993443	601.5512	14817.1191	202404
## 3	3	Slope_CO2	0.9999907	168.3200	10075.2183	202404
## 4	4	Slope_CH4_low	0.9996407	633.0314	-320.6965	202404
## 5	5	Slope_CH4_high	0.9993443	601.5512	14817.1191	202404
## 6	6	Slope_CO2	0.9999907	168.3200	10075.2183	202404



0.2 Now calculate the CH4 & CO2 concentrations in ppm

```
#head(raw)

#pull out methane standards
Samples <- raw %>%
  filter(!str_detect(Sample_Type, "Standard")) %>%
  filter(!str_detect(Sample_Type, "STD_CO2")) %>%
  filter(!str_detect(Sample_Type, "Blank")) %>%
  filter(!str_detect(Sample_Type, "Lab air")) %>%
  filter(!str_detect(Sample_Type, "ChkStd")) %>%
  filter(!str_detect(Sample_ID, "Shakey")) %>%
  filter(!str_detect(Sample_Type, "NA"))
#head(Samples)

#Now flag any areas that are above the 100ppm area for CH4
Samples$CH4_Curve <- ifelse(Samples$CH4_Area > 71000, "High", "Low")
#head(Samples)

#Calculate CH4 concentrations in ppm
Samples$CH4_Conc_ppm <- ifelse(Samples$CH4_Area > 71000, (Samples$CH4_Area - Slope_CH4_high$Intercept) / Slope_CH4_high$Slope, 0)

#Calculate CO2 concentrations in ppm
#Samples$CO2_Conc_ppm <- ((Samples$CO2_Area - Slope_CO2$Intercept) / Slope_CO2$Slope)
```

```

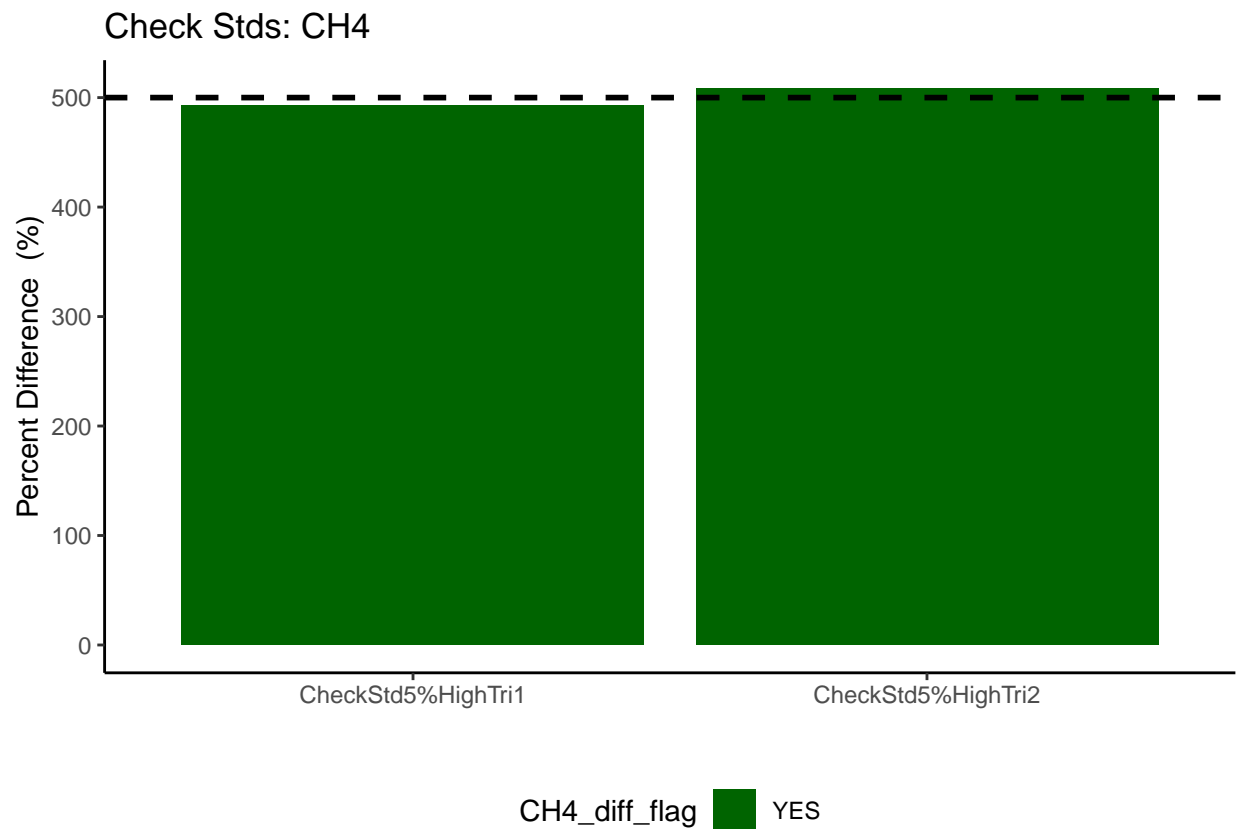
#head(Samples)

#####make flags for any dilutions needed
#highest CH4 standard = 10000
#highest CO2 standard = 50000

Samples$CH4_Flag <- ifelse(Samples$CH4_Conc_ppm >10000, "Needs Dilution", "Within Range")
#Samples$CO2_Flag <- ifelse(Samples$CO2_Conc_ppm >50000, "Needs Dilution", "Within Range")
#head(Samples)

```

0.3 Check the Check Standards



0.4 Dilution correct samples

```

#multiply the concentration by the dilution factor
Samples$CH4_Conc_ppm_dilcorr <- (Samples$CH4_Conc_ppm * Samples$Dilution_Factor)

#Samples$CO2_Conc_ppm_dilcorr <- (Samples$CO2_Conc_ppm * Samples$Dilution_Factor)

#check results
#head(Samples)

```

```

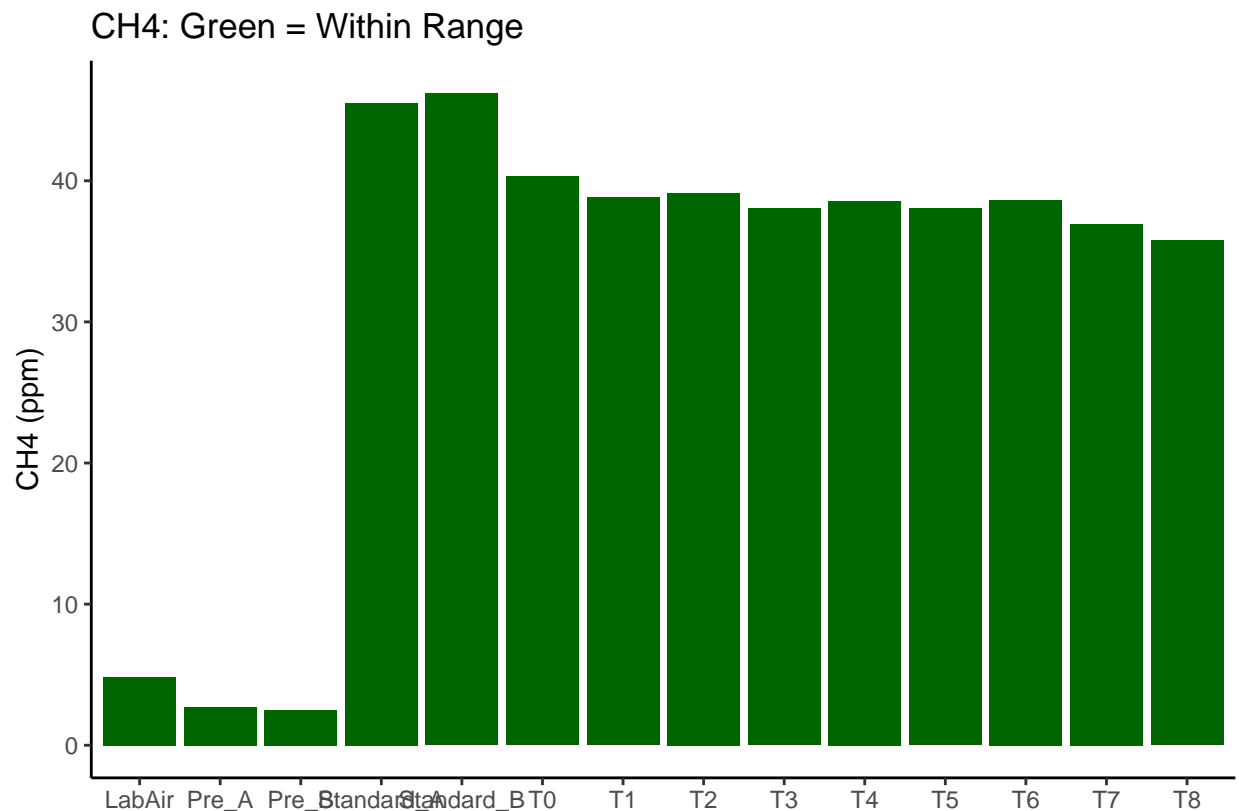
#quick first look at the samples
ch4_samples <- ggplot(data = Samples, aes(x = Sample_ID, y = CH4_Conc_ppm, fill=CH4_Flag)) +
  geom_bar(stat = 'identity') +
  scale_fill_manual(values=c("darkgreen","red"))+
  #scale_fill_gradient2(low='red', mid='white', high='blue', space='Lab') +
  theme_classic() + labs(x= " ", y="CH4 (ppm)", title="CH4: Green = Within Range") +
  theme(legend.position="none")

#co2_samples <- ggplot(data = Samples, aes(x = Sample_ID, y = CO2_Conc_ppm, fill=CO2_Flag)) +
#  geom_bar(stat = 'identity') +
#  scale_fill_manual(values=c("red", "darkgreen"))+
#  #scale_fill_gradient2(low='red', mid='white', high='blue', space='Lab') +
#  theme_classic() + labs(x= " ", y="CO2 (ppm)", title="CO2: Green = Within Range") +
#  theme(legend.position="none")

#ggarrange(ch4_samples, co2_samples, nrow=1, ncol=2)

print(ch4_samples)

```



0.5 If samples are water calculate gas in water

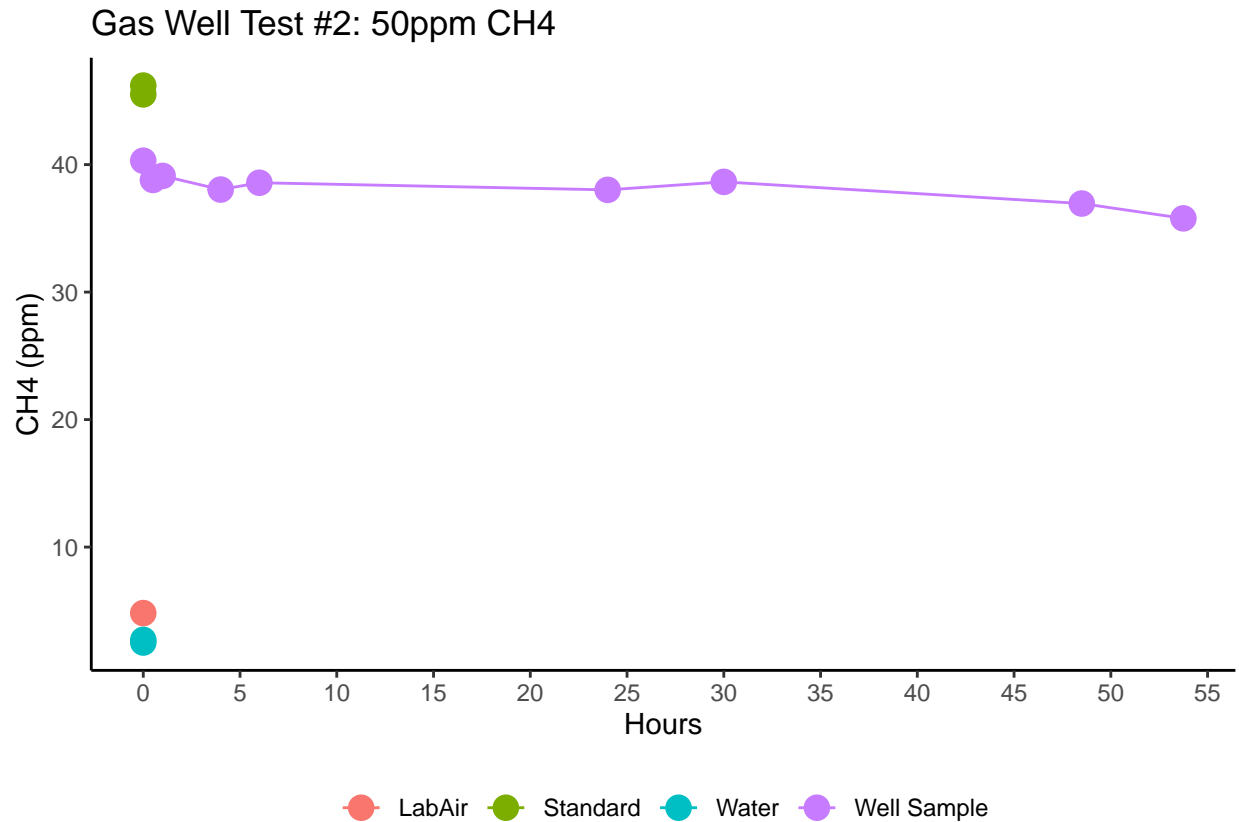
0.6 Write out processed data & slopes

```
#check results
#head(Samples)

#pull out what we need
#Samples1 <- Samples[ ,c(1:3,6:17,21:24)]
#head(Samples1)

final_data <- Samples %>%
  #select(Project, Plot, grid, sample_name, Vial_ID, date, ) %>%
  mutate(
    Project = "COMPASS",    # new column with same value on every row
    Experiment = "TEMPEST: Gas Well Test #2")#,
    #Run_notes = run_notes    # new column with notes about the run
  #)

#plot the data thru time so we can see the equilibration
final <- ggplot(data = final_data, aes(x = Hours, y = CH4_Conc_ppm_dilcorr, color=Event_Stamp)) +
  geom_point(size=4) +
  geom_line() +
  scale_x_continuous(breaks = seq(0, 55, by = 5))+
  theme_classic() + labs(x= "Hours", y="CH4 (ppm)", title="Gas Well Test #2: 50ppm CH4") +
  theme(legend.position = "bottom", legend.title=element_blank())
final
```



```
#this needs altered to match the tempest metadata and clean up
final_data <- final_data %>%
  rename(
    CH4_ppm = CH4_Conc_ppm_dilcorr ,
    CH4_uM = CH4_Conc_umol,
    # add more rename pairs as needed
  ) %>%
  select(
    Project, Experiment, Sample_Year, Sample_Month, Sample_Day, Sample_Time,
    Sample_ID, Event_Stamp, Hours,
    CH4_ppm, CH4_uM, CH4_Flag #, tdn_mgL, tdn_uM, tdn_flag, Analysis_runtime,
    #Run_notes
    # list columns in the order you want them
  )

head(final_data)
```

```
##   Project                Experiment Sample_Year Sample_Month Sample_Day
## 1 COMPASS TEMPEST: Gas Well Test #2      2025   September      16
## 2 COMPASS TEMPEST: Gas Well Test #2      2025   September      16
## 3 COMPASS TEMPEST: Gas Well Test #2      2025   September      16
## 4 COMPASS TEMPEST: Gas Well Test #2      2025   September      16
## 5 COMPASS TEMPEST: Gas Well Test #2      2025   September      16
## 6 COMPASS TEMPEST: Gas Well Test #2      2025   September      16
##   Sample_Time Sample_ID Event_Stamp Hours   CH4_ppm   CH4_uM   CH4_Flag
## 1         1200   LabAir   LabAir      0  4.822081 0.003941833 Within Range
```


## 2	1000	Pre_A	Water	0	2.718543	0.111114294	Within Range
## 3	1000	Pre_B	Water	0	2.513579	0.102736843	Within Range
## 4	1000	Standard_A	Standard	0	45.505453	0.037198644	Within Range
## 5	1000	Standard_B	Standard	0	46.197922	0.037764706	Within Range
## 6	1000	TO Well Sample		0	40.303769	0.032946503	Within Range

```
write.csv(final_data, "Processed Data/TMP_GasWellTest_2_GHG_Processed.csv")
```

```
#end
```