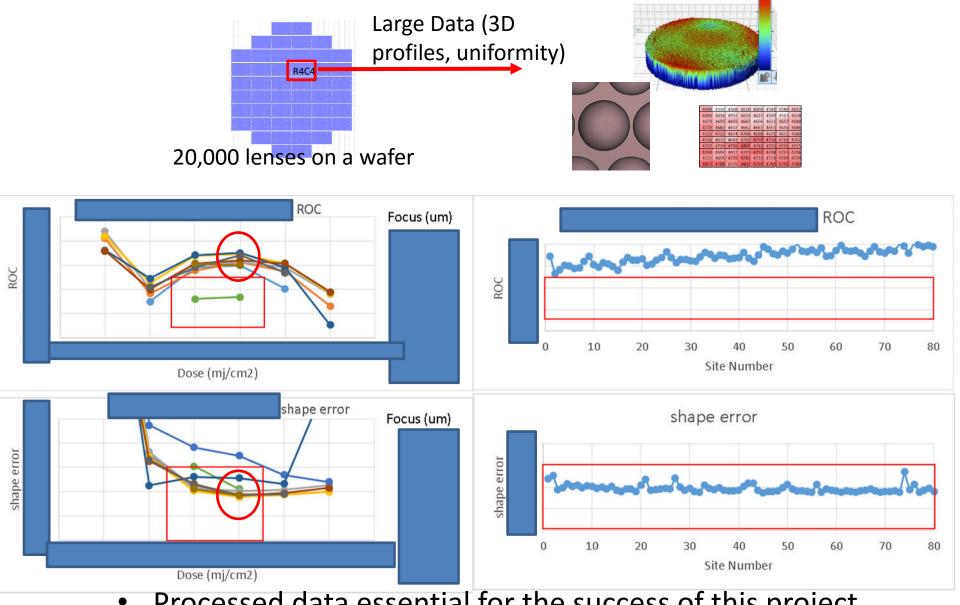
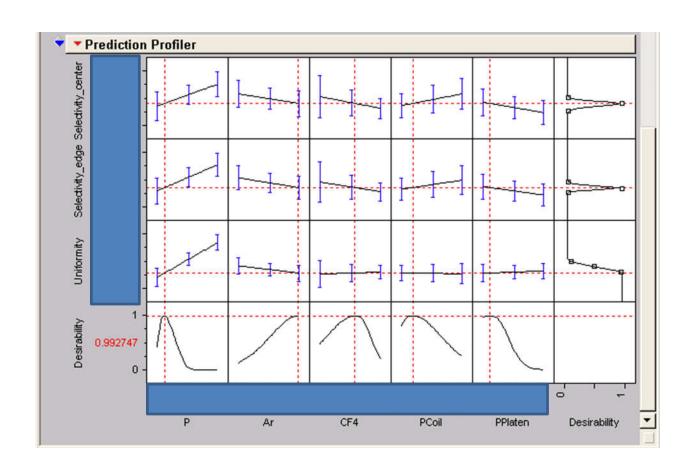
Microlens Array (IMT, 2013-present)



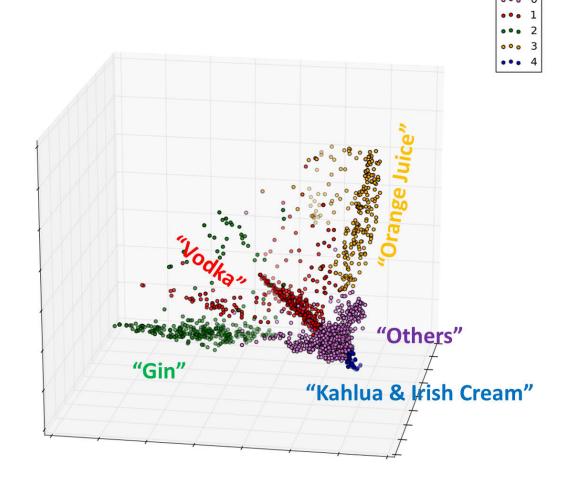
- Processed data essential for the success of this project.
- Transitioning to a volume production phase

### JMP (IMT, 2013-present)



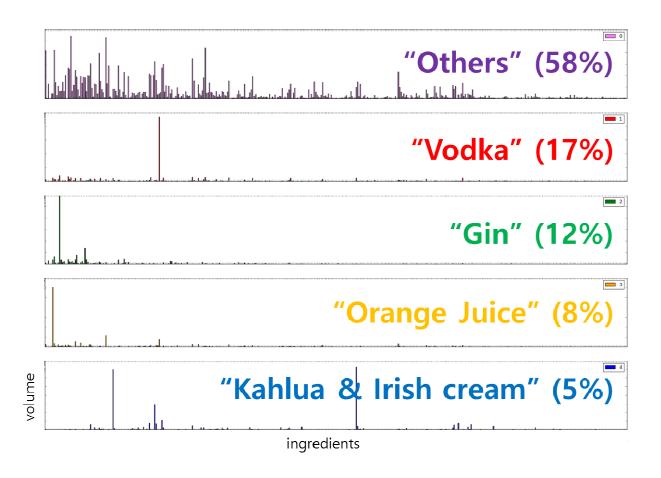
SAS JMP used for optimizing process

### Python programming as a hobby (2016)



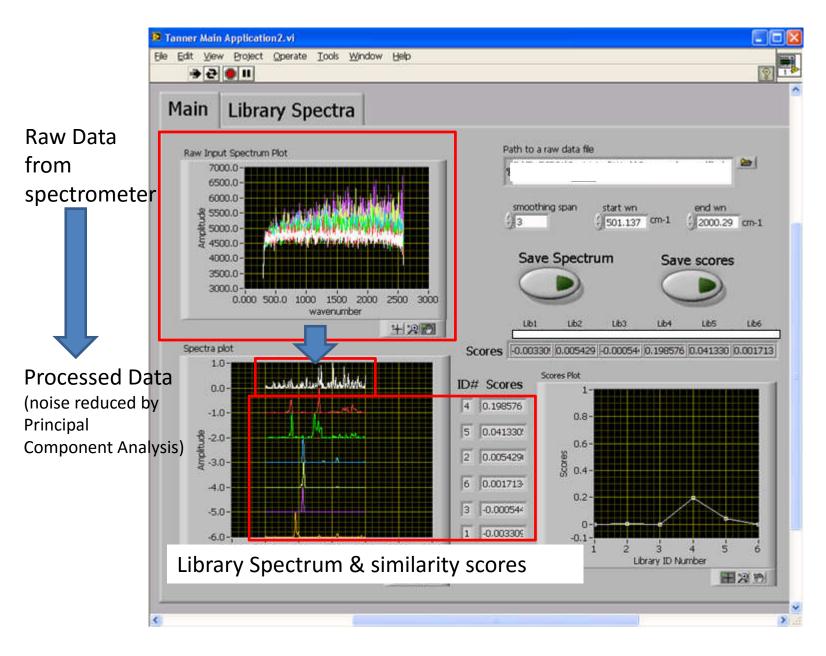
- 3081 cocktail recipes were downloaded from <a href="http://www.TheCocktailDB.com">http://www.TheCocktailDB.com</a> using their API.
- Each cocktail was embedded in a 471-dimensional vector space using ingredient information.
- Principal Component Analysis (PCA) was used in order to visualize 3081 cocktails in a 3D space.
- Spectral Cluster algorithm was used to separate the groups
- Generated profiles of each market segment.

### Python programming as a hobby (2016)



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### Chemical Detection (Tanner Research, 2009-2013)



### **Chemical Detection (Tanner Research, 2009-2013)**

Name	Formula for Scores (worst=0, best=1)
Euclidean Distance (ED)	$S_k = \frac{L_k \cdot U}{\sqrt{L_k \cdot L_k} \sqrt{U \cdot U}}$ (=cos $\theta_k$ , where $\theta_k$ is the angle between $L_k$ and U)
Correlation (CO)	$S_k = \frac{{L_k}' \cdot U'}{\sqrt{{L_k}' \cdot {L_k}'} \sqrt{U' \cdot U'}}$ , where $L_k' = L_k - \frac{\sum_{i=1}^n L_{ki}}{n}$ and $U' = U - \frac{\sum_{i=1}^n U_i}{n}$
Absolute Value (AV)	$S_k = 1 - \frac{\sum_{i=1}^{n}  L_{ki} - U_i }{n}$
Least Squares (LS)	$S_k = 1 - \frac{\sum_{i=1}^n (L_{ki} - U_i)^2}{n}$
Single Chemical Library Regression (PCR)	$S = UL^t(LL^t)^{-1}$ , assuming $U = SL$ holds true.

n: number of spectral data points (e.g., 1024)

U: unknown spectrum vector (e.g., 1x1024)

L: library spectra matrix (e.g., 6x1024)

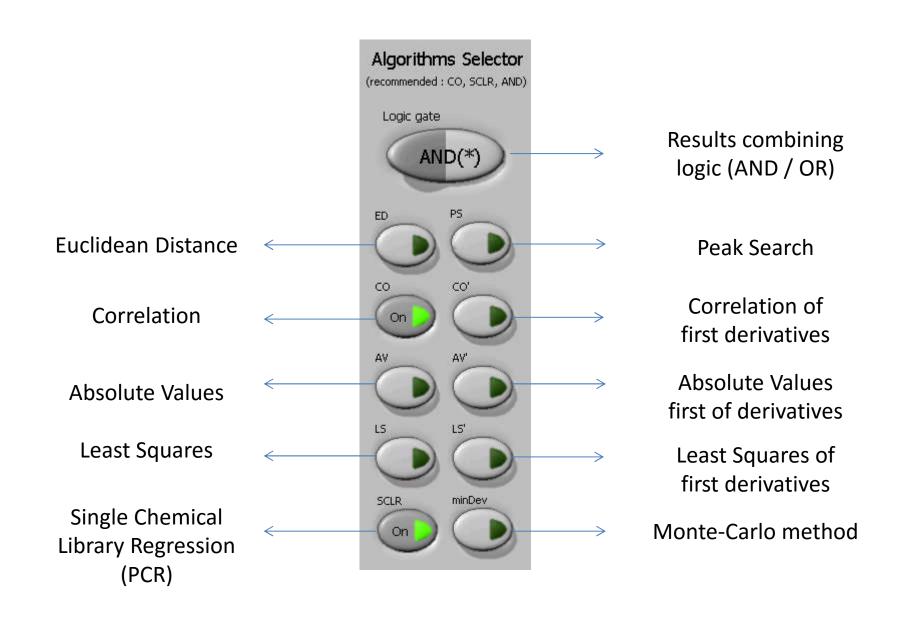
S: score matrix (e.g., 1x6)

 $L_k$ : k-th library spectrum vector (e.g., 1x1024)

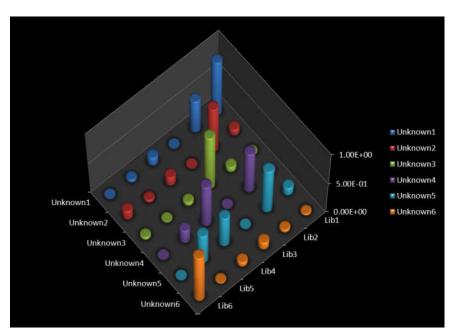
 $L_{ki}$ : i-th spectral component of k-th library spectrum vector

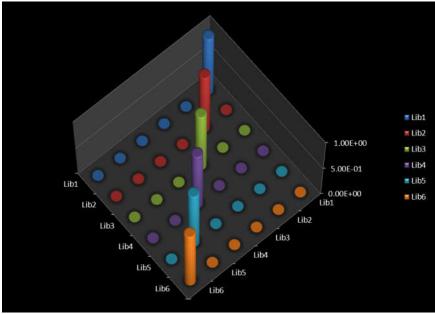
 $S_k$ : score of unknown vector against k-th library spectrum vector

#### **Chemical Detection (Tanner Research, 2009-2013)**

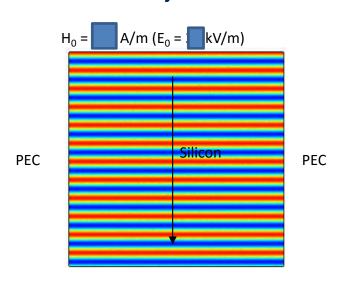


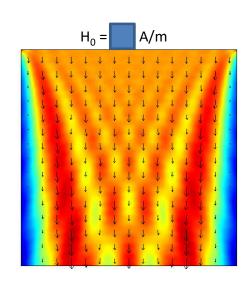
### **Chemical Detection Algorithms**

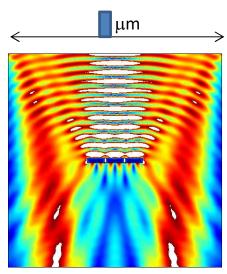


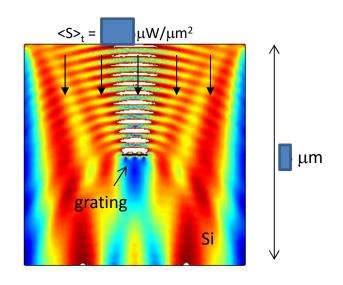


 This standoff-chemical detection system passed rigorous field tests (accuracy, recall) at customer's proving ground. Customer later asked us to deliver software only.



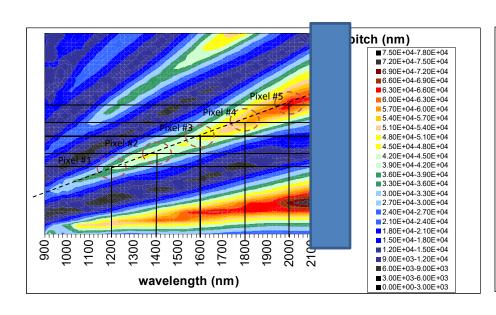


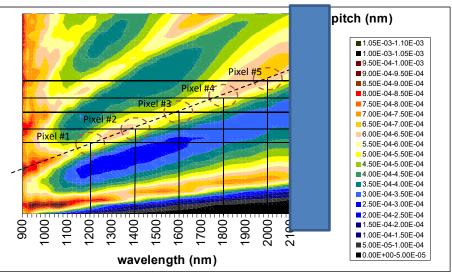




Strong E-fields and large  $\Delta T$  when pitch =  $(2n-1)/2*\lambda_{spp}$ , n = 1,2,3, ... Weak E-fields and small  $\Delta T$  when pitch =  $n*\lambda_{spp}$ , n = 1,2,3, ... Reference: D. Pacifici, et al., PRB 77, 115411 (2008)

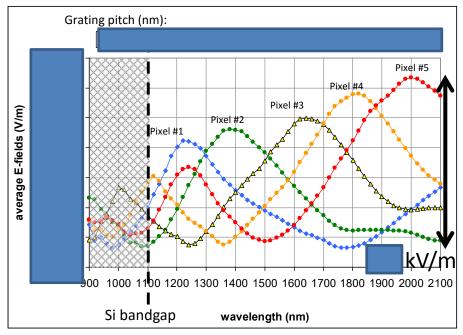
E-fields  $\Delta T$ 

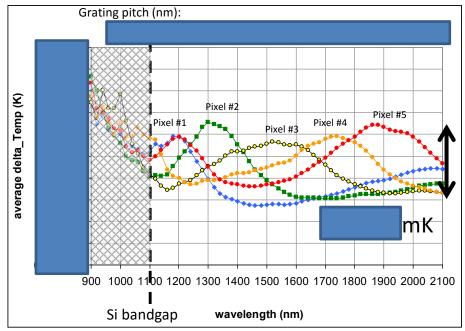


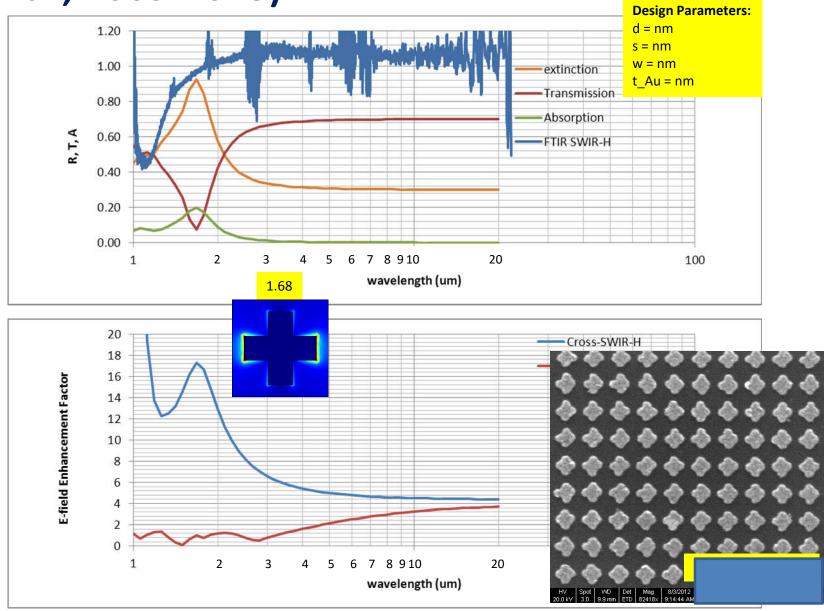


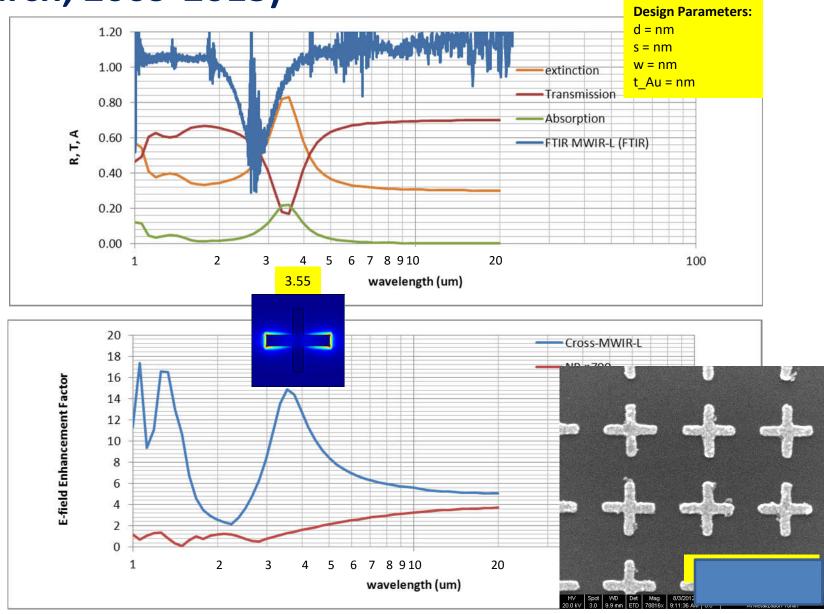
Average E-fields on the gaps

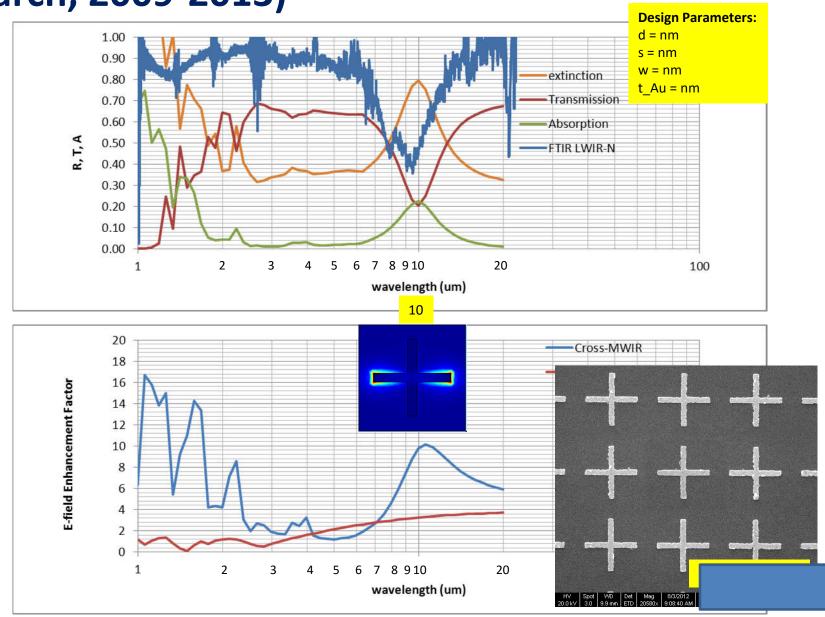
average  $\Delta T$  (temperature change due to ohmic heating) on the gaps

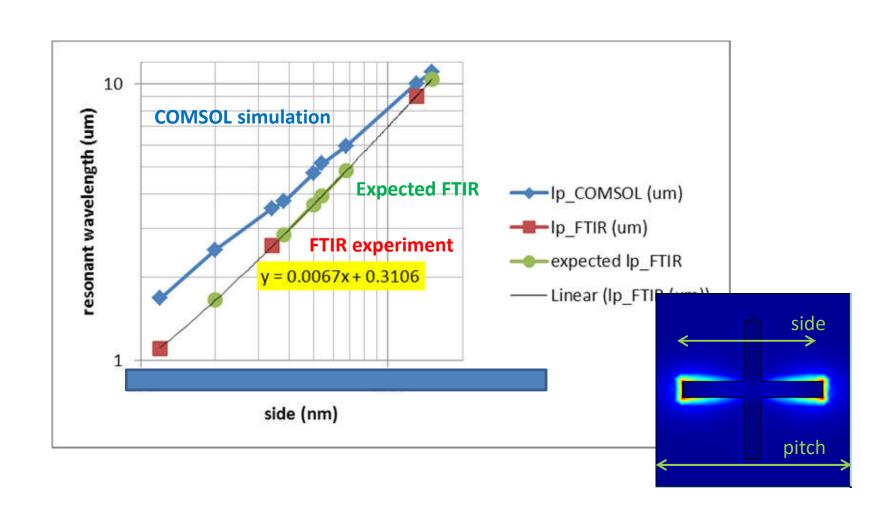




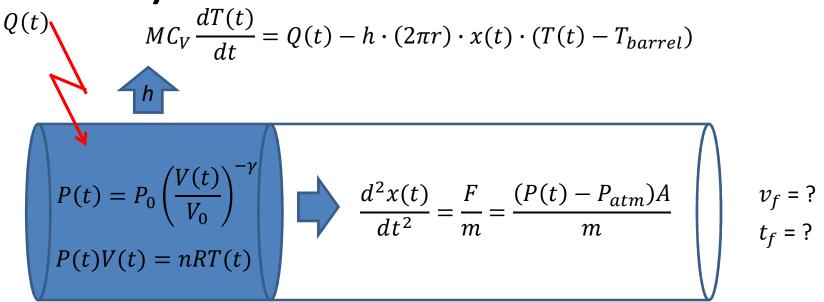


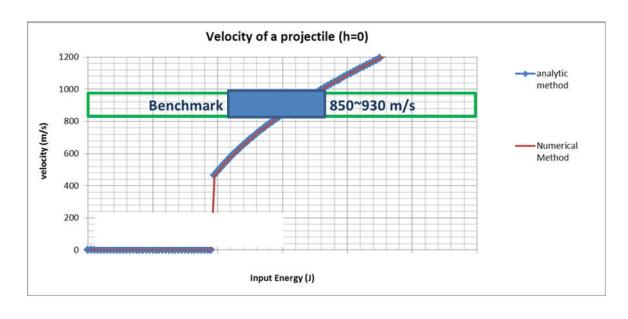


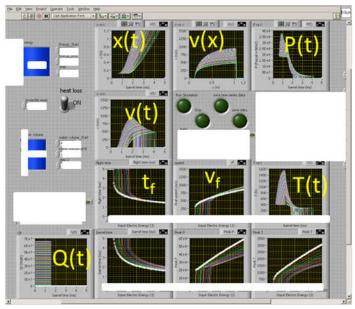




## Numerical Methods Simulation (Tanner Research, 2009-2013)

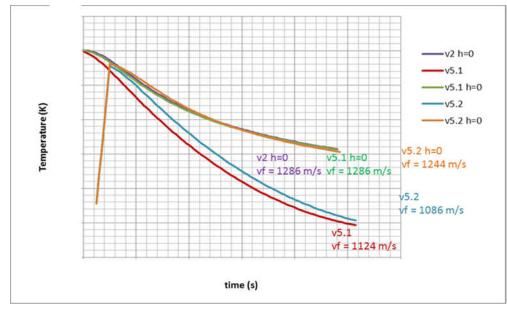


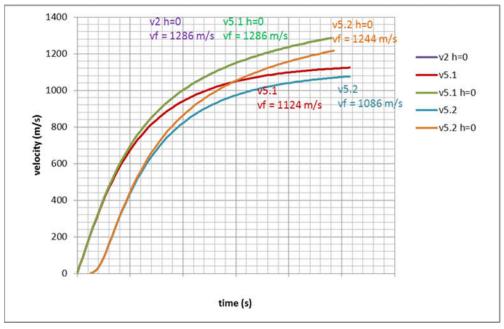




## Numerical Methods Simulation (Tanner Research,

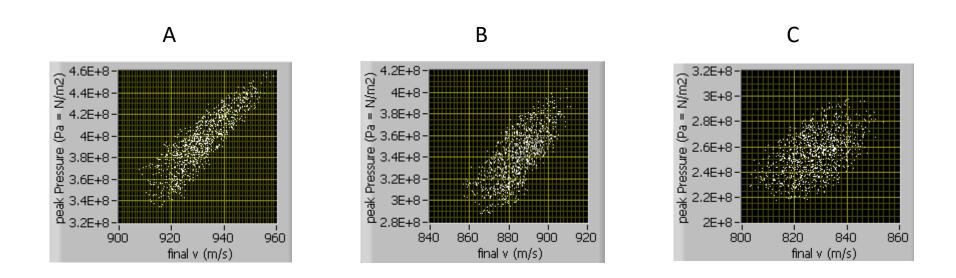
2009-2013)





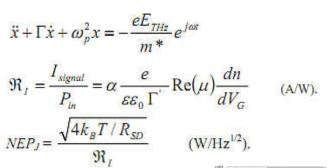
# Numerical Methods Simulation (Tanner Research, 2009-2013)

#### Monte-Carlo Method (optimization)

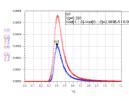


- Peak pressure related to the breakage of the system.
- Final v relevant to the performance of the system.

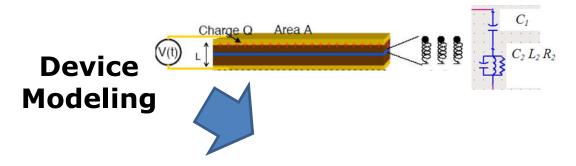
### THz Detector (UCSB Ph.D. 2003-2009)



Device Model Revision



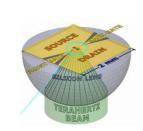




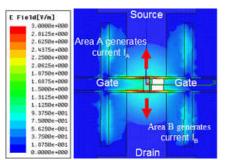


#### **Fabrication**

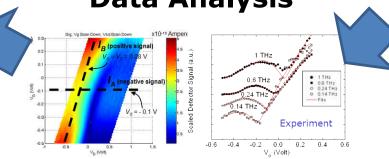




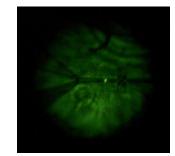
## FEM EM Simulation



#### **Data Analysis**

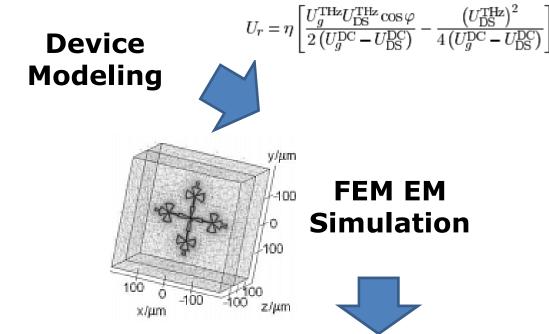


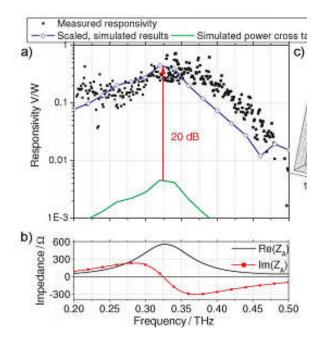
#### **Measurements**



### THz Mixer (Tanner Research, 2009-2013)

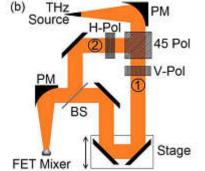
 $j_{SD} = n(\omega_G)ev(\omega_{SD})$ 





**Data Analysis** 

### Measurements



### **Cleanroom Fabrication**



## Thank you!