# TODOp

- 1. Find how garrett made charts in his manuscript
- 2. Make charts better or change input data
- 3. Do comparisons of Cl species with F species
  - (a) Garrett seems to have started doing that in lagrangian hi res mass frac.ipynb, but he doesn't include measured Cl data.

### cfs2canteraPD

- 1. Get injected\_species from netj\_file
- 2. Get conc\_cutoff from netj\_file (inject conc)
- 3. Make it run PFRTP-Garrett and cfs\_post if needed

### $cfs\_post$

- 1. Make trajectory selection faster
- 2. Make it run PFRTP-Garrett if needed

#### PFRTP-Garrett

- 1. Make interactive widget to choose run conditions
- 2. Fix thing with flame temperature estimate
- 3. Temperatures shouldn't go down then back up

### Instructions

- 1. Run netj\_generate.ipynb to generate input files from lists of run conditions
- 2. Run cfs\_batchrun.ipynb to send the generated .netj files to CFS and run the simulations
- 3. Run PFRTP-Garrett.ipynb to use Cantera to perform the same simulations.
- 4. Run cfs\_post.ipynb, and select .netj files.
- 5. Run cfs2canteraPD.ipynb to generate path diagrams using Cantera.

## **Problems**

- 1. There are many many .ipynb files in Modeling\_GPD, many of which are near duplicates of each other, scattered across folders. It is difficult to tell which ones are the most up to date and which ones are obsolete.
- 2. Residence Time column from cantera has no duplicate times. CFS will have 3-7 rows of identical times
- 3. Cantera increases time steps constantly, CFS has variable time steps
- 4. Concentrations of non-major species in cantera are around 1e-100, in CFS they are around 1e-15
  - (a) Produces problems for reaction path diagrams because cantera can't tell which is the major species
  - (b) Concentrations are all 0 until injection, then they all instantly increase
    - i. This is probably because the transient solution solves from the injection point and then tries to work backwards to the flame. Injecting at the flame may fix this.
- 5. CF3Or has 15 reactions, CF2s has 74, CF4 15, C2F6 6. C2F6 behavior in path diagrams may be caused by incomplete reactions
  - (a) C2F6 may just not have that many reactions because it is a larger molecule, and the C-C bond is the most likely to break
- 6. Several reactions are "ambiguous" and cantera will use the designated default reaction
  - (a) Default reaction may not be conducive for simulating  ${\rm C2F6}$
- $7.\ \,$  Some reactions are not parsed (they may just not be relevant)

# Notes

- 1. CF4 will never be destroyed at temps below 1200K, 100% DE for temps 1600 C
- 2. Higher values of MAXIT seem to produce more residence time steps

Table 1: Files, Folders, and Descriptions

File	Folder	Description
${\rm flow\_convert.ipynb}$	PFAS_Modeling/	Converts liters per minute to kilograms per second for air and for CH4 using cantera
netj_generate.ipynb	${\rm PFAS\_Modeling}/$	Takes a set of varying run conditions and produces netj files to do all iterations
$cfs\_batchrun.ipynb$	${\it PFAS\_Modeling/}$	Uses CFS batch method to run generated netj files, prompts user for inputs to determine simulation duration
PFRTP-Garrett.ipynb	PFAS_Modeling/Pseudo_PFR	Uses a PFR simulation in Cantera to make similar results as CFS
$cfs\_post.ipynb$	${\it PFAS\_Modeling/}$	Runs CFS in GUI mode to automate post-processing of vtk files to include species concentration
cfs2canteraPD.ipynb	${\it PFAS\_Modeling/}$	Converts vtk to csv data, then produces a path diagram using Cantera for a specified trajectory, residence time, and threshold.
pfr.ipynb	${\rm PFAS\_Modeling}/$	Example PFR that came with Cantera
pfr2.ipynb	PFAS_Modeling/	Example Cantera PFR that includes heat loss
vtk2csv.py	PFAS_Modeling/	Stand alone script for converting a specific vtk to a csv
lagrangian hi res mass frac.ipynb	L:/Lab/AMCD_PFAS_Incineration/Modeling_GPD/ Cantera/PFR DE S Curves/	Makes plots of multiple species DE% vs Temperature
eulerian <injectant>.ipynb</injectant>	L:/Lab/AMCD_PFAS_Incineration/Modeling_GPD/ Cantera/influent concentration/	Make multidimensional color plots for conc. vs Temperature vs DE%. Also absolute and relative yield vs conc.
paths < SR >. ipynb	$\begin{array}{c} L:/Lab/AMCD\_PFAS\_Incineration/Modeling\_GPD/\\ Cantera/paths/ \end{array}$	Makes path diagrams for several species at specified SR
T99.ipynb	$\begin{array}{c} L:/Lab/AMCD\_PFAS\_Incineration/Modeling\_GPD/\\ Cantera/T99/ \end{array}$	Makes T99 plots of multiple species
lagrangian.ipynb	L:/Lab/AMCD_PFAS_Incineration/Modeling_GPD/ Cantera/PFR DE S Curves/archive/	Makes plots of species DE vs Temperature. Plots from this script don't look the same as plots in the doc, so it is probably outdated.

Table 2: Validated Path Diagrams - CFS

Table 21 Validated Latin Blagfalling CLS													
		CHF3				C2F6							
		Por	Ports				Ports						
	1	4	6	8	1	4	6	8	1	4	6	8	
27.5 kW									×			×	
45 kW	×	<b>√</b>							×	×		×	

Table 3: Validated Path Diagrams - Cantera

		C1	F4		CHF3				C2F6				
		Ports				Ports							
	1	4	6	8	1	4	6	8	1	4	6	8	
27.5 kW	×	×	×	×					<b>√</b>	<b>√</b>	<b>√</b>	<b>√</b>	
45 kW	<b>√</b>	<b>√</b>	<b>√</b>	<b>√</b>					<b>√</b>	<b>√</b>	<b>√</b>	<b>√</b>	

Table 4: CFS Executed Simulations

		CF	<b>'</b> 4			СН	F3		C2F6				
	Ports					Por	rts		Ports				
	1	4	6	8	1	4	6	8	1	4	6	8	
27.5 kW		<b>√</b>							<b>√</b>	<b>√</b>	<b>√</b>	<b>√</b>	
45 kW	<b>√</b>	✓			<b>√</b>	<b>√</b>	<b>√</b>		✓	✓	<b>√</b>	<b>√</b>	

Table 5: Cantera Executed Simulations

			C	F4			CH	IF3		C2F6				
		Ports				Ports				Ports				
ľ		1	4	6	8	1	4	6	8	1	4	6	8	
	$27.5~\mathrm{kW}$	<b>√</b>	<b>√</b>	<b>√</b>	<b>√</b>					<b>√</b>	<b>√</b>	<b>√</b>	<b>√</b>	
ľ	$45~\mathrm{kW}$	<b>√</b>	<b>√</b>	<b>√</b>	<b>√</b>					<b>√</b>	<b>√</b>	<b>√</b>	<b>√</b>	