## Additional Functions for PFAS Suspect List

## **Function Description**

For the analysis and useful application of the suspect list, additional functions were generated for R to process the dataset.

## Set up R environment

To open Microsoft Excel files, you must install the 'xlsx' package.

```
install.packages("xlsx")
```

get\_suspectlist: To download the most recent version of the PFAS suspect list, use this function.

```
source("fn/get_suspectlist.R")
get_suspectlist("PFAS Suspect List.xlsx")
# You can designate the list be saved as a different file
# name if you want.
```

open\_suspectlist: Open the suspect list into the R environment (you must do this before any other steps)

```
source("fn/open_suspectlist.R")
suspectlist <- open_suspectlist("PFAS Suspect List.xlsx")
## use the file location if this is different</pre>
```

add\_ionization: calculates the exact m/z value for indicated ionization states.

add\_residual: calculates the residual mass after removing the repeating units (default = CF2)

expand\_additional: all delimited information stored in the ADDITIONAL column is expanded as additional columns. NOTE: There will be a significant number of new columns with NA values for most compounds.

```
source("fn/expand_additional.R")
suspectlist_w_additional <- expand_additional(suspectlist)
write.csv(suspectlist_w_additional, "suspectlist_w_additional.csv",
    row.names = FALSE)</pre>
```

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