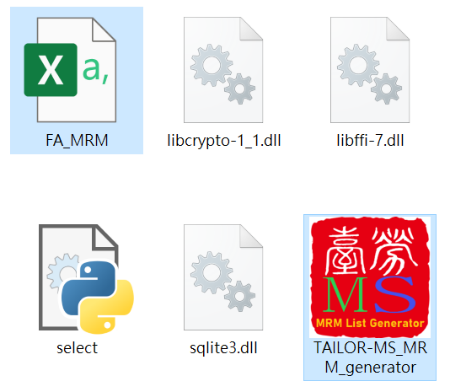
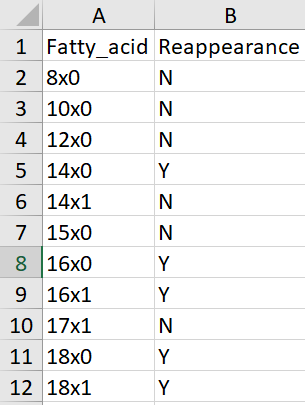
**TAILOR-MS**

**TAILOR-MS** (TriAcylglycerol Identifier for LOw Resolution Mass Spectrometry) is a Python package for automating the triacylglycerol identification process with input targeted LC-MS (or other types of mass spectrometric data) data. It consists of two independent scripts: **MRM list generator** and **Identifier**.

**MRM List Generator**

****This script creates an MRM list for setting up targeted mass spectrometric acquisition methods for TG with [M+NH4]+/[ M+NH4-RCOONH4]+ precursor/product ion pairs. After unzipping TAILOR-MS\_MRM\_generatior.zip file, you will find in folder “**FA\_MRM.csv**” and “**TAILOR-MS\_MRM\_generator.exe**” files. To use MRM generator, simply replace contents in FA\_MRM.csv and double-click to run the exe file.

**Figure 1. Partial contents of unzipped TAILOR-MS\_MRM\_generator folder→**

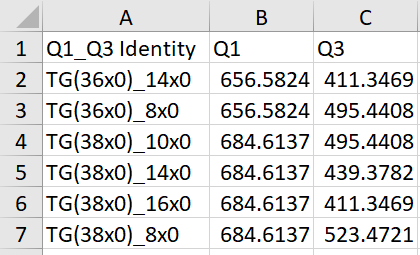
*Input file*

FA\_MRM.csv must contain the following contents:

**Fatty\_acid:** fatty acyl groups one intends to use for generating possible TG structures. “x” instead of “:” are used here so that Microsoft Excel does not automatically convert the input FAs to time.

**Reappearance:** indicates if the user allows an FA group to appear 2 or 3 times in a TG structure when generating MRM lists *eg* TG(18:1\_18:1\_18:1) and TG(8:0\_18:1\_18:1) will be allowed here but not TG(8:0\_8:0\_8:0).

**←Figure 2. An example for FA\_MRM.csv**

*Output files*

TAILOR-MS MRM generator creates two files: **MRM\_list.csv** and **TG\_by\_MRM.csv**.

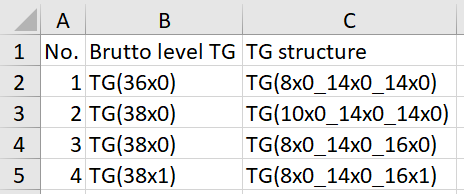
Below are the contents of **MRM\_list.csv:**

**Q1\_Q3 Identity:** brutto level TGs and neutral loss FAs corresponding to Q1/Q3 pairs

**Q1:** Q1 m/z from LipidMaps

**Q3:** Q3 m/z from LipidMaps

**↑Figure 3. An example of MRM\_list.csv**

TG\_by\_MRM.csv contains the following contents:

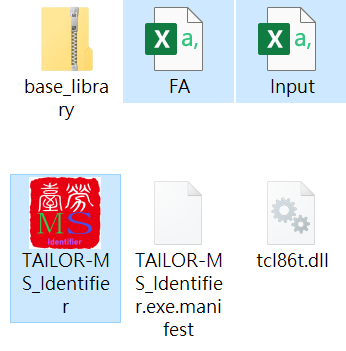
**No.:** numbers

**Brutto level TG:** brutto level TG

**TG structure:** structures with FA details

**↑Figure 4. An example of TG\_by\_MRM.csv**

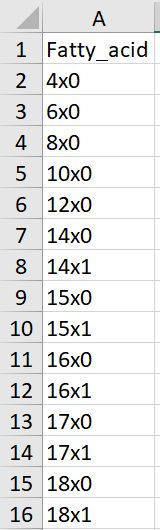
**Identifier**

TAILOR-MS Identifier identifies and predicts TG structures with fatty acyl information but without positional information. It reads two input files: **“FA.csv”** and **“Input.csv”**. To use TAILOR-MS Identifier, simply replace the contents of these two csv files with your own data and preferred settings and double click **TAILOR-MS\_Identifier.exe** to run the script.

First, unzip TAILOR-MS\_Identifier.zip file and find the following files in the folder. Part of the folder should look like this:

**Figure 5. Partial contents of unzipped TAILOR-MS\_Identifier folder→**

**↓Figure 6. An example of FA.csv file**

FA.csv must contain a single column with FA groups to be used for structural identification and prediction, similar to the first column of “FA\_MRM.csv” file for MRM list generator.

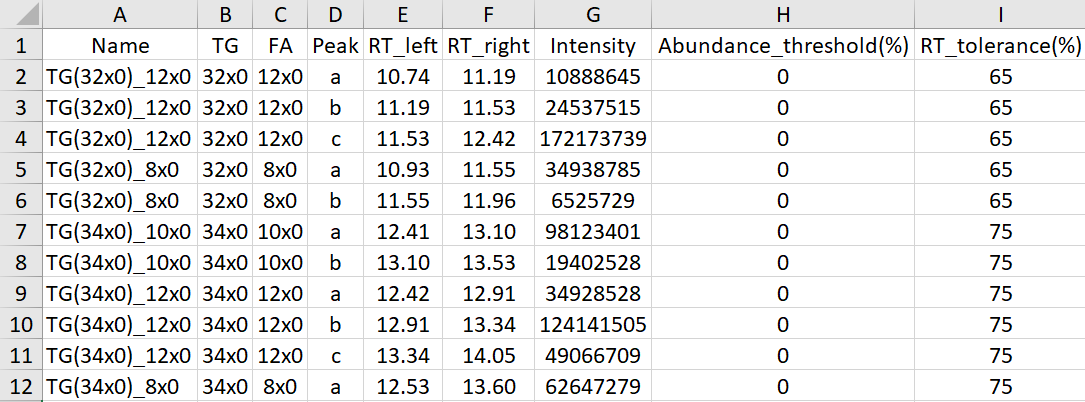
Input.csv file has several columns and are explained below**:**

**Name:** anything the user wants to name their peaks. Must be English characters and numbers.

**TG:** brutto level TG.

**FA:** FA neutral losses.

**Peak:** used to assign peaks captured under the same Q1/Q3 which indicates chromatographic separation of structural isomers. Must be lower case English alphabets (a-z).

**↑Figure 7. An example of Input.csv file**

**RT\_left:** left boarder of peak retention time. Must be 0 or positive values.

**RT\_right:** right boarder of peak retention time. Must be positive values and greater than corresponding RT\_left values.

**Intensity:** peak intensities, which can be areas, heights or concentrations. Must be positive values.

**Abundance\_threshold(%):** a user-defined threshold. Any structures with a relative abundance (vs the peak with largest intensity of all peaks having the same TG brutto level) lower than this threshold will be removed. Must be 0 or positive values. All peaks with the same brutto level TG must have the same set value.

**RT\_tolerance(%):** a user defined threshold. To confirm the existence of a TG structure, the constituting peaks of this structure must overlap. RT\_tolerance enables the user to exclude TG structures of which the peaks do not overlap to the % above the set values. More detailed explanation can be found in TAILOR-MS paper. Must be positive values. All peaks with the same brutto level TG must have the same set value.

*Output file:*

Double click TAILOR-MS\_Identifier.exe to start the run. Depending on the size of input data, this may take a few minutes.

**“Results.csv”** will be spawned in the same folder. Contents of this file are explained bellow.

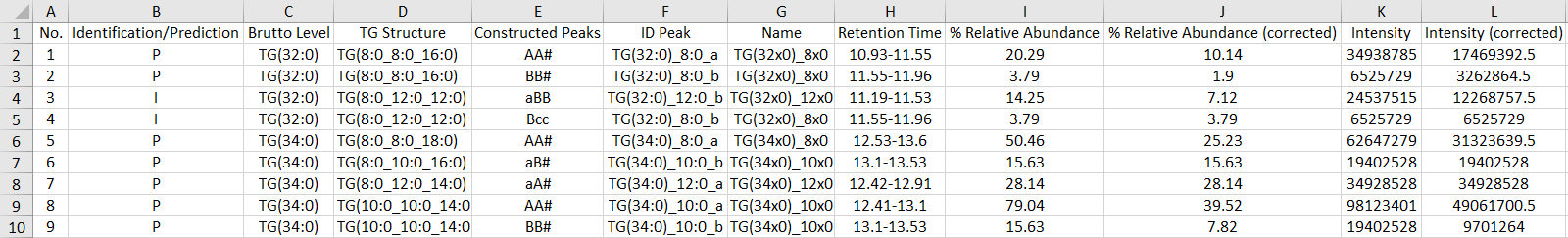
**No.:** numbers

**Identification/Prediction:** indicate if the TG structure has been identified (having 3 FA groups in acquisition method and match the rules) or predicted (having only 2 FA groups)

**Brutto Level**: brutto level TG

**TG Structure:** TG structures with FA details

**Constructed Peaks:** correspond to “Peak” information from Input.csv. Capitalised letter(s) indicates the ID peak; # indicates missing peak from the acquisition method (predicted TG structures).

**ID Peak:** the FA peak with minimal intensity among the three FA groups that constitute a specific TG, which can be used to indicate the abundance of this particular TG structure.

**Name:** names from Input.csv.

**Retention Time:** retention time spans of ID peak.

**% Relative Abundance:** uncorrected relative abundances of ID peak.

**% Relative Abundance (corrected):** relative abundances corrected when ID peak reoccurs in a structure.

**Intensity:** uncorrected intensities of ID peak.

**Intensity (corrected):** intensities corrected when ID peak reoccurs in a structure.

**←Figure 8. An example of Results.csv**