

Act as a professional researcher in organic chemistry. You are asked to extract information from a given `reaction_text` and export the information to a ORD JSON record. A JSON record is an ORD JSON if it uses the Open Reaction Database (ORD) schema.

I will describe the 5 steps of this task. I will then guide you step by step to perform this task on one exemplar `reaction_text`. I will then give you a new `reaction_text` and you need to generate the JSON record.

Step 1: Identify all the chemicals in the given `reaction_text`. An chemical identifier can be the name of a compound, for example, `methanol`. An identifier can also be an index or a generic description, for example, `compound 6`, or `desired compound`.

Step 2: <CONTINUE TO DEFINE STEPS>

Here is the first example. `reaction_text` is the text between two delimiters `` and ``. The exported ORD JSON record is the text between two delimiters ### and ###.

`reaction_text` = ``<EXAMPLE REACTION TEXT>``

Here is the workflow how to extract information from this `reaction_text` and export them to a ORD JSON record.

Step 1: This reaction involves six chemicals. 4-aminobenzotriazole, acetic acid, potassium bromide, ammonium molybdate, hydrogen peroxide, desired product.

Step 2: <CONTINUE TO DEFINE STEPS>

example_ORD_JSON = ###<EXAMPLE ORD JSON>###

Here is the second example.

<CONTINUE TO THE SECOND EXAMPLE>