

## The Basic Trial Solution Method

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The method is outlined here for a second order differential equation  $ay'' + by' + cy = f(x)$ . The method applies unchanged for  $n$ th order equations.

- Step 1. Repeatedly differentiate the atoms of  $f(x)$  until no new atoms appear. Collect the distinct atoms so found into a maximal list of  $k$  atoms. Multiply these atoms by **undetermined coefficients**  $d_1, d_2, \dots, d_k$ , then add, defining **trial solution**  $y$ .
- Step 2. Substitute  $y$  into the differential equation.

**Fixup Rule I.** If some variable  $d_p$  is missing in the substituted equation, then step 2 fails. Correct the trial solution as follows. Variable  $d_p$  appears in trial solution  $y$  as term  $d_p A$ , where  $A$  is an atom. Multiply  $A$  and all its related atoms  $B$  by  $x$ . The modified expression  $y$  is called a **corrected trial solution**. Repeat step 2 until the substituted equation contains all of the variables  $d_1, \dots, d_k$ .

- Step 3. Match coefficients of atoms left and right to write out linear algebraic equations for  $d_1, d_2, \dots, d_k$ . Solve the equations for the unique solution.
- Step 4. The corrected trial solution  $y$  with evaluated coefficients  $d_1, d_2, \dots, d_k$  becomes the particular solution  $y_p$ .

## Symbols

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The symbols  $c_1, c_2$  are reserved for use as arbitrary constants in the general solution  $y_h$  of the homogeneous equation. Symbols  $d_1, d_2, d_3, \dots$  are reserved for use in the trial solution  $y$  of the non-homogeneous equation. Abbreviations:  $c = \text{constant}$ ,  $d = \text{determined}$ .

## Superposition

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The relation  $y = y_h + y_p$  suggests solving  $ay'' + by' + cy = f(x)$  in two stages:

- (a) Apply the linear equation **recipe** to find  $y_h$ .
- (b) Apply **the basic trial solution method** to find  $y_p$ .

We expect to find two arbitrary constants  $c_1, c_2$  in the solution  $y_h$ , but in contrast, no arbitrary constants appear in  $y_p$ . Calling  $d_1, d_2, d_3, \dots$  *undetermined* coefficients is misleading, because in fact they are eventually *determined*.

## Fixup rule II

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The rule predicts the corrected trial solution  $y$  without having to substitute  $y$  into the differential equation.

- Write down  $y_h$ , the general solution of homogeneous equation  $ay'' + by' + cy = 0$ , having arbitrary constants  $c_1, c_2$ . Create the corrected trial solution  $y$  iteratively, as follows.
- Cycle through each term  $d_p A$ , where  $A$  is an atom. If  $A$  is also an atom appearing in  $y_h$ , then multiply  $d_p A$  and each **related atom** term  $d_q B$  by  $x$ . Other terms appearing in  $y$  are unchanged.
- Repeat until each term  $d_p A$  has atom  $A$  distinct from all atoms appearing in homogeneous solution  $y_h$ . The modified expression  $y$  is called the **corrected trial solution**.

### Fixup rule III

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The rule predicts the corrected trial solution  $\mathbf{y}$  without substituting it into the differential equation. This iterative algebraic method uses the atom list of the homogeneous equation to create  $\mathbf{y}$ .

- Write down the roots of the characteristic equation. Let  $L$  denote the list of distinct atoms for these roots.
- Cycle through each term  $d_p A$ , where  $A$  is a atom. If  $A$  appears in list  $L$ , then multiply  $d_p A$  and each **related atom** term  $d_q B$  by  $x$ . Other terms appearing in  $\mathbf{y}$  are unchanged.
- Repeat until the atom  $A$  in an arbitrary term  $d_p A$  of  $\mathbf{y}$  does not appear in list  $L$ .<sup>a</sup> The modified expression  $\mathbf{y}$  is called the **corrected trial solution**.

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<sup>a</sup>The number  $s$  of repeats for initial term  $d_p A$  equals the multiplicity of the root  $r$  which created atom  $A$  in list  $L$ .

### Fixup rule IV

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The rule predicts the corrected trial solution  $\mathbf{y}$  without substituting it into the differential equation. This algebraic method uses the roots of the characteristic equation to create  $\mathbf{y}$ .

- Write down the roots of the characteristic equation as a list  $\mathbf{R}$ , according to multiplicity.
- Let  $\mathbf{G}$  denote a largest group of related atom terms in  $\mathbf{y}$  with first atom  $\mathbf{A}$ . If  $\mathbf{R}$  contains a root  $r$  of multiplicity  $s$ , and an atom  $\mathbf{B}$  for  $r$  is related to atom  $\mathbf{A}$ , then multiply all terms of  $\mathbf{G}$  by  $x^s$ . If no root in  $\mathbf{R}$  has atom related to  $\mathbf{A}$ , then no action is taken.
- Repeat the previous step for all groups of related atoms in  $\mathbf{y}$ . The modified expression  $\mathbf{y}$  is called the **corrected trial solution**.

