

The Corrected Trial Solution
in
the Method of Undetermined Coefficients

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Definition of Related Atoms

Atoms A and B are *related* if and only if their successive derivatives share a common atom. Then x^3 is related to x and x^{101} , while x is unrelated to e^x , xe^x and $x \sin x$. Atoms $x \sin x$ and $x^3 \cos x$ are related, but the atoms $\cos 2x$ and $\sin x$ are unrelated.

The Basic Trial Solution Method

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.** Extract all distinct atoms from $f(x)$, $f'(x)$, $f''(x)$, ... to construct 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.

Basic Correction Rule. 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

The symbols $\mathbf{c}_1, \mathbf{c}_2$ are reserved for use as arbitrary constants in the general solution \mathbf{y}_h of the homogeneous equation.

Symbols $\mathbf{d}_1, \mathbf{d}_2, \mathbf{d}_3, \dots$ are reserved for use in the trial solution \mathbf{y} of the non-homogeneous equation. Abbreviations: $\mathbf{c} = \text{constant}$, $\mathbf{d} = \text{determined}$.

Superposition

The relation $\mathbf{y} = \mathbf{y}_h + \mathbf{y}_p$ suggests solving $a\mathbf{y}'' + b\mathbf{y}' + c\mathbf{y} = \mathbf{f}(x)$ in two stages:

- (a) Find \mathbf{y}_h as a linear combination of atoms computed by applying Euler's theorem to factors of $ar^2 + br + c$.
- (b) Apply the **basic trial solution method** to find \mathbf{y}_p .
 - We expect to find two arbitrary constants c_1, c_2 in the solution \mathbf{y}_h , but in contrast, no arbitrary constants appear in \mathbf{y}_p .
 - Calling d_1, d_2, d_3, \dots *undetermined* coefficients is misleading, because in fact they are eventually *determined*.

Annihilator Polynomial $q(r)$ for $f(x)$

Assume $f(x)$ is a linear combination of atoms. Each atom corresponds to, according to Euler's Theorem, a factor $(r - a - ib)^{k+1}$ of a polynomial. Choose the largest k , over all atoms sharing this factor, and then multiply all maximal factors, taken over all atoms appearing in $f(x)$, to obtain a polynomial $q(r)$. If a complex factor appears in $q(r)$, then multiply $q(r)$ by the conjugate of that factor. Finally, $q(r)$ is real. It is called an *annihilator polynomial* for $f(x)$.

Example: If $f(x) = x + x^2 + \cos x$, then the maximal factors according to Euler's Theorem are r^3 and $(r - i)$. Then $q(r) = r^3(r - i)(r + i)$, because complex factors must appear with their conjugates. Finally, $q(r) = r^3(r^2 + 1)$.

Annihilator Differential Equation $Aw = 0$ for $f(x)$

Assume $f(x)$ is a linear combination of atoms and $q(r)$ is an annihilator polynomial for $f(x)$. Then $q(r)$ is the characteristic polynomial of a higher order linear homogeneous equation $Aw = 0$ and $w = f(x)$ is a particular solution of this differential equation. The equation $Aw = 0$ is called an *annihilator differential equation* for $f(x)$.

Example: If $f(x) = x + x^2 + \cos x$, then $q(r) = r^3(r^2 + 1) = r^5 + r^3$ and $Aw = 0$ is the differential equation $w^{(5)} + w^{(3)} = 0$. We verify that $f(x)$ satisfies

$$\begin{aligned} A(f) &= f^{(5)} + f^{(3)} \\ &= ((x + x^2)^{(5)} + (x + x^2)^{(3)}) + ((\cos x)^{(5)} + (\cos x)^{(3)}) \\ &= 0 + 0 \\ &= 0. \end{aligned}$$

Correction rule I: The Annihilator Method

The rule computes the corrected trial solution \mathbf{y} without having to substitute \mathbf{y} into the non-homogeneous differential equation $L\mathbf{y} = \mathbf{f}$.

- Let $p(r)$ be the characteristic polynomial for the homogeneous differential equation $L\mathbf{y} = \mathbf{0}$, from which we obtain the homogeneous general solution $\mathbf{y}_h(x)$.
- Let $q(r)$ be an annihilator polynomial for $\mathbf{f}(x)$ and $A\mathbf{w} = \mathbf{0}$ its annihilator differential equation, so that $A(\mathbf{f}) = \mathbf{0}$. *We never need to find $A\mathbf{w} = \mathbf{0}$ explicitly!*
- Multiply $p(r)$ and $q(r)$ to obtain $p(r)q(r)$, which is the characteristic equation of $A(L\mathbf{y}) = \mathbf{0}$. Then $\mathbf{y}(x)$ is a solution of $A(L\mathbf{y}) = \mathbf{0}$, because $A(L\mathbf{y}) = A(\mathbf{f}) = \mathbf{0}$. Expand \mathbf{y} as a linear combination of atoms selected from the maximal factors of $p(r)q(r)$.
- The superposition principle $\mathbf{y}(x) = \mathbf{y}_h(x) + \mathbf{y}_p(x)$ implies the corrected trial solution $\mathbf{y}(x)$ is obtained by removal of all atoms shared with $\mathbf{y}_h(x)$.

Correction rule II

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**.

Correction rule III

The rule predicts the corrected trial solution y without substituting it into the differential equation. This iterative algebraic method uses the atom list of the homogeneous equation to create 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 y are unchanged.
- Repeat until the atom A in an arbitrary term $d_p A$ of y does not appear in list L .^a The modified expression y is called the **corrected trial solution**.

^aThe number s of repeats for initial term $d_p A$ equals the multiplicity of the root r which created atom A in list L .

Definition of function **atomRoot**

- $\text{atomRoot}(x^j e^{rx}) = r$ for r real.
- $\text{atomRoot}(x^j e^{ax} \cos bx) = \text{atomRoot}(x^j e^{ax} \sin bx) = a + ib$.

Correction rule IV

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 correct \mathbf{y} .

- Write down the roots of the characteristic equation as a list \mathbf{R} , according to multiplicity.
- Subdivide trial solution \mathbf{y} into groups \mathbf{G} of related atoms, by collecting terms and inserting parentheses.
- If a group \mathbf{G} contains an atom \mathbf{A} with $r = \text{atomRoot}(\mathbf{A})$ in list \mathbf{R} , then multiply all terms of \mathbf{G} by x^s , where s is the multiplicity of root r .
- Repeat the previous step for all groups \mathbf{G} in \mathbf{y} . The modified expression \mathbf{y} is called the **corrected trial solution**.

