Hydrogen Bonding Network Emulating Frequency Driven Source of Triangular Pulses

Rostislav Rusev, George Angelov, Elitsa Gieva, Marin Hristov, and Tihomir Takov

Abstract—A three output microelectronic circuit functionally equivalent to a hydrogen bonding network is modeled. Proton transfer characteristics of each hydrogen bond of the network are emulated by block-elements in the microelectronic circuit with their respective I-V characteristics. These characteristics are coded in Matlab where the dynamic and static analyses are carried out. The results imply that in static mode the functionally analogous circuit operates as a current source or an amplifier. In dynamic mode the circuit behaves as a voltage driven triangular pulse signal source. The simulations show that the generated pulses at the three circuit outputs have different frequency, amplitude, and width.

Index Terms—hydrogen bonding network, proton transfer, microelectronic circuit.

I. INTRODUCTION

The application of molecular interactions such as hydrogen bonding for design of molecular self-assemble systems in micro- and nanoelectronics, including liquid crystals [1], molecular solids [2], polymeric assemblies [3] has attracted the attention of scientists long time ago. For example, linear porphyrine layers are self-organized through hydrogen bonding coordinated the self-organization of the lipid bi-layer membranes. The electron acceptor is placed on one side of the membrane and the electron donor is placed on the other side. Photo-current arises when the sample is irradiated by white light. The photo-current can be observed only between ordered structures and it cannot be observed between separate molecules. In this way, a photonic electron conductor from simple molecule components can be constructed. On the other hand, the development of such new systems is increasing because the hydrogen bonding in anisotropic molecular self-assembled structures can form bridges between synthetic materials and/or bio-functional systems. In the biofunctional nature systems, e.g. proteins, the hydrogen bonding is crucial for molecule recognition, molecule stabilization, dynamic reversibility [4], etc. Proton transfer proteins such as Bacteriorhodopsin (bR), which transfers proton current via its own hydrogen bonding network (HBN), might be used for interesting applications in microelectronics. Its proton current versus voltage characteristic is similar to the output characteristic of a microelectronic circuit (differential amplifier) [5].

Similarly to Bacteriorhodopsin we are investigating the possibility for information transfer via branched HBN of β-lactamase protein.

II. MODEL AND EQUATION

The object of the present investigation is a microelectronic circuit similar to β-lactamase branched HBN [6]. The proton transfer parameters of each hydrogen bond are also investigated in [6]. These parameters are the basis of the presented microelectronic simulations in this paper. Here the K-El.pot characteristics of the network (K – proton transfer parameter and El.pot – electrostatic potential) are interpreted as I-V characteristics where the proton transfer parameter is regarded as the electrical current.

The probability for proton transfer from the donor to acceptor is proportional to the value of the K parameter. The proton current between donor and acceptor of each hydrogen bond depends on the K parameter as well. Some authors [7] consider the K parameter to be equivalent to the proton current.

We calculate the proton transfer parameter (K) using the following equations:

\[
K = \frac{k_B T}{2\pi} \exp\left(-\frac{E_b - h\omega / 2}{k_B T}\right) \tag{1}
\]

where: \(K\) – proton transfer parameter, \(k_B\) – Boltzmann constant, \(E_b\) – energy barrier, \(h\) – Planck constant, \(\omega\) – frequency, \(T\) – temperature [K].

The energy barrier is calculated by:

\[
E_b = (s_A (R(DA) - t_A)^2 + v_A) + s_B E_{12} + (s_C \exp(-t_C (R(DA) - 2)) + v_C) (E_{12})^2 \tag{2}
\]

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The branched hydrogen bonding network is represented in Fig. 1 and its functionally analogous electric circuit is shown in Fig. 4.

The $K-V$ characteristic of each bond is given in Fig. 2 and Fig. 3. It can be observed that the $K-V$ characteristics are nonlinear. This allows for the hydrogen bonds to process the signals similarly to the semiconductor devices. On this basis we can construct three- and four-terminal microelectronic block elements that emulate the operation of the hydrogen bonds.

The block elements are based on heavy atoms that are part of the hydrogen bonding network. Each heavy atom simultaneously appears as donor and acceptor in the hydrogen bonding network. In this reason the input of each block element corresponds to the acceptor part of the heavy atom and its output corresponds to the donor part of the heavy atom.

The equivalent circuit composed of these microelectronic block-elements is given in Fig. 4 where each block-element corresponds to a heavy atom in the hydrogen bonding network. The arginine residue R164 is depicted as the four terminal block-element T1. Since it is strong proton donor in the network of hydrogen bonds, in the analogous electric circuit it is represented as voltage-driven current source (with three outputs). Output currents and voltages of $T_1$ are different and they are driven by the input voltage. The second block-element $T_2$ is analogous to aspartic acid residue $D_179$, which is strong proton acceptor. In the electric circuit the block-element $T_2$ is the first output.

Block-elements $T_4$ and $T_5$ are analogous to the input molecule $w_{295}$ and aspartic acid residue $D_{176}$ of Fig. 1. They are shown as three-terminal block-elements. Input and output voltages of each of the $T_4$ and $T_5$ block-element are equal and their input but output currents are different. On the other hand $T_5$ is the second circuit output.

The block-element $T_7$ of the circuit is analogous to the strong proton acceptor of HBNE171. It forms two hydrogen bonds with its oxygen $OE_1$ and $OE_2$ atoms: at one of its sides is $R_{164} NH_1...(OE_1)_{171}E$ and at the other side is $w_{859} OH...(OE_2)_{171}E$. The potential on $OE_2$ atom as well as proton current between $w_{859}OH$ and $(OE_2)_{171}E$ both depend on the potential of $OE_1$ because of cooperative effect. Consequently, the output current $I_7$ and voltage $U_7$ of the analogous block-element $T_7$ depend on the input voltage $U_{71}$. 

The $K-V$ characteristics of hydrogen bonds: OD2 is oxygen atom of Aspartic acid residue D176, NH1 is nitrogen of Arginine residue R164, OE1 is oxygen atom of Glutamic acid residue E171, and OH is oxygen atom of water molecules $w_{295}$, $w_{753}$, $w_{859}$.

The block elements are based on heavy atoms that are part of the hydrogen bonding network. Each heavy atom simultaneously appears as donor and acceptor in the hydrogen bonding network. In this reason the input of each block element corresponds to the acceptor part of the heavy atom and its output corresponds to the donor part of the heavy atom.
The block-element T8 appears to be the last output of the circuit that is analogous to w859 of the HBN. We have to point out that the next water molecule of the HBN, w753, is taken into account but it is not shown in the electrical circuit because the input current to w753 remains the same regardless of how many water molecules are bound after it (under the assumptions of macroscopic model). In this reason we included only block-element T8. Similarly to the other water molecule w295 it is described with three-terminal block-element with identical input and output voltages and different currents.

The relations between currents and voltages of the block-elements in the microelectronic circuit are given by polynomials.

Equation (4) below describes the first output of T1:

\[ U_1 \in [-1.3:3.2] \text{ step 0.1} \]  

\[ I_1 = -5.10^{-5} (U_1)^3 - 8.10^{-5} (U_1)^2 - 2.10^{-5} (U_1) + 0.0071 \]  

Next are the equations of T2

\[ U_2 = 0.9994 (U_1) - 0.3421 \]  

\[ I_2 = I_1 \]  

The equation of the second output of T1:

\[ U_3 = 1.0066 (U_1) - 0.0967 \]  

\[ I_3 = 0.0005 (U_3)^3 - 0.0006 (U_3)^2 - 0.0006 (U_3) + 0.028 \]  

Equations for T4

\[ U_4 = 0.9802 (U_3) + 0.4294 \]  

\[ I_4 = -0.1922 (U_4)^3 + 0.2821 (U_4)^2 + 0.0044 (U_4) + 16.922 \]  

Equations for T5 which is the second output of the circuit

\[ U_5 = 0.9374 (U_4) + 1.4074 \]  

\[ I_5 = I_4 \]  

Equations of the third output of T1

\[ U_6 = 1.0193 (U_1) + 0.3216 \]  

\[ I_6 = -0.0072(U_6)^3 + 0.0069 (U_6)^2 - 0.03 (U_6) + 1.2646 \]  

Equations of block-element T7

\[ U_7 = 1.0387 (U_6) - 0.5498 \]  

\[ U_7 = 0.9705 (U_71) - 0.5167 \]  

\[ I_7 = 0.0591 (U_7)^3 + 0.0162 (U_7)^2 - 0.6792 (U_7) + 4.5597 \]  

Equations of T8 that is the third output of the electrical circuit

\[ U_8 = -0.0437 (U_7)^2 + 1.0318 (U_7) + 0.4173 \]  

\[ I_8 = 0.0041 (U_8)^4 - 0.0456 (U_8)^3 + 0.0772 (U_8)^2 + 0.234 (U_8) + 1.8401 \]  

These equations are coded in Matlab where the static and dynamic analyses of the microelectronic circuit are performed [8]. Sample Matlab code is given in Fig. 5.

### III. STATIC ANALYSIS

The static analysis is performed by feeding input voltage \( U_{in} \) between \(-1.3 \) and \(+3.2 \text{ [V]} \). Subsequently, a comparison between I-V characteristics of each block-element and the proton transfer characteristics of the respective hydrogen bond from paper [6] is carried out. Sample results of the juxtaposition are depicted on Fig. 6 and 10; for the other block-elements the results are similar and they are not shown on the figures. As it can be seen, the polynomials well describe the hydrogen bond functions. The maximal standard deviation from all simulations versus data set is \( \pm 4.33 \% \). The model is in good agreement with the experimental data for hydrogen bonds because standard deviation of the measurements is in the range of \( 4\% \) to \( 12\% \).
Still, the simulations show that all output voltages are proportional to the input voltage (Fig. 7). Uout1 and Uout3 are almost equal and Uout2 is shifted up. When Uin is varied between $-1.3$ and $+3.2$ [V] then Uout1 and Uout3 are changing in interval $-1.5 \div +3.2$ [V], and Uout2 is increasing from 0.5 to 4.5 [V].

In contrast to the voltages, the currents are always positives. The currents of the different outputs are changing in different intervals from $10^{-3}$ to $10^1$ [pA]. The I-V characteristics of output 1 and output 2 have similar forms (Figs. 8 and 9) but they are shifted to one another. For output 1: when Uout1 is in the interval between $-1.5$ and 0 [V] Iout1 is constant.

Consequently, this circuit can be used as current source. When Uout1 is varying between 0 and +3 [V] the current is decreasing, i.e. this output is similar to the output of an amplifier. By analogy such conclusions can be made also for output 2: the circuit operates as current source at Uout2 = +0.5 to +2.5 [V] because the Iout2 current is stable. When Uout2 is between +2.5 to +4.5 [V], the output of the circuit is again similar to an amplifier output.
IV. Dynamic Analysis

The dynamic analysis is done by transient simulations in the range $0 \div 10^{-10}$ [s] because in nature objects proton transfers durations are in these time intervals. Subsequently the analogous microelectronic circuit should operate at much higher frequencies — ca. 10 [GHz]. The circuit is tested with sine input voltages with different amplitudes and frequencies. At $U_{in} = 1.5 \times \sin(t \times 5 \times 10^{11})$ the outputs have characteristics (not shown in the paper) and the circuit operates as signal modulator producing signals with different amplitude and form.

At $U_{in} = -1 + \sin(t \times 1.5 \times 10^{11})$ stable circuit operation is observed (Fig. 11): at the three outputs triangular pulses are generated (Figs. 12, 13, 14).

Figures show that the pulses are with equal frequency, phase, and positive amplitude. The three pulses are with different amplitudes and the width of the pulses at the first output is less than the width of the pulses at outputs 2 and 3. On the other hand we can control the period of the output pulses by changing the period of the input voltage. For instance, on Fig. 15 is given characteristic of $I_{out2}$ versus time that is obtained at $U_{in} = -1 + \sin(t \times 5 \times 10^{11})$, analogous to the characteristics obtained at the other outputs. By comparing Figs. 13 and 15 it can be seen that the triangular pulses of Fig. 15 are with larger frequency. Similar result is observed for the other outputs and thence they are not shown.

Finally, the microelectronic circuit that is analogous to network of hydrogen bonds can operate as frequency driven source of triangular pulses. The three outputs generate pulses with different amplitude and width.
V. CONCLUSION

The hydrogen bonding network model proved that such biofunctional system has real microelectronics applications. The analogous microelectronic circuit may operate in static mode as a current source or an amplifier. In dynamic mode the microelectronic circuit may function as a voltage driven triangular pulse signal generator. At the three circuit outputs might be generated pulses with different frequency, amplitude, and width. Thus, the biocircuit appears to be extremely flexible and is applicable to multiple circuit purposes.

REFERENCES


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