

Table 1:

<i>Design</i>	<i>Thickness (nm)</i>	$E_3 - E_2$ (meV)	$E_2 - E_1$ (meV)	f_{21} (a.u.)	$\Delta_{1'3}$ (meV)
<i>D1</i>	4.4/ <u>5.6</u> / 1.2 /7.7/ 2.8 /8.1	34.8	16.3	0.32	6.7
<i>D2</i>	4.4/ <u>5.6</u> / 1.2 /7.7/ 2.4 /7.9	34.7	16.5	0.45	6.6
<i>D3</i>	4.2/ <u>5.5</u> / 1.2 /8.1/ 2.0 /7.8	34.6	16.4	0.62	6.7
<i>D4</i>	4.0/ <u>5.5</u> / 1.2 /8.5/ 1.7 /7.5	34.3	16.3	0.82	6.9

TABLE I. Overview of the layer sequences and key parameters of the four designs. The layer thicknesses of one active module are given with the $\text{Al}_{0.15}\text{Ga}_{0.85}\text{As}$ barriers in bold and GaAs wells in plain text. The underlined wells are doped with Si at the level of $2.66 \times 10^{16} \text{ cm}^{-3}$, $2.61 \times 10^{16} \text{ cm}^{-3}$, $2.62 \times 10^{16} \text{ cm}^{-3}$ and $2.58 \times 10^{16} \text{ cm}^{-3}$ for D1 – D4 respectively, yielding average bulk level of $5.0 \times 10^{15} \text{ cm}^{-3}$. $E_3 - E_2$ is the energy separation between subbands 3 and 2, $E_2 - E_1$ is the energy separation between subbands 2 and 1, f_{21} is the scaled oscillator strength, and $\Delta_{1'3}$ is the anticrossing gap energy of the subbands 1' and 3. The principal difference between the four designs is the optical transition diagonality, which is characterized by f_{21} .