

## Rate-Independent Constructs for Chemical Computation

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### Appendix: Logarithm Reactions

We present chemical reactions that implement the pseudo-code presented in the text.

**System Initialization** As with our other modules, we have a small set of reactions to control the overall timing of our system.



**Halving** We use a slight variation of our decrement module to implement the operation  $x = x / 2$ .



**Increment** We use our increment module to implement the line  $y = y + 1$ . We set  $y$  to be 1 initially.



**Cleanup** Once the module has completed, we decrement  $y$  by one, storing the result in  $y_f$ .

$$x'_{ab} + y'_{ab} \xrightarrow{\text{slow}} done^P \quad (17)$$

$$done^P + 2x \xrightarrow{\text{fast}} 2x \quad (18)$$

$$done^P + x' \xrightarrow{\text{fast}} x' \quad (19)$$

$$done^P + y' \xrightarrow{\text{fast}} y' \quad (20)$$

$$done^P \xrightarrow{\text{slow}} done \quad (21)$$

$$2done \xrightarrow{\text{fast}} done \quad (22)$$

$$done + 2y \xrightarrow{\text{slow}} y + y_f \quad (23)$$

**Absence Indicators** Two special absence indicators are used by the halving and increment modules above; a total of 13 are needed for the system to function properly. They are of the same form as all other absence indicators, described in the paper.