```
[1] import tellurium as te
      te.newTiledFigure(rows=2, cols=1)
      <tellurium.plotting.engine_plotly.PlotlyTiledFigure at 0x7f28fc4125f8>
👼[2] // Created by libAntimony v2.9.3
      model *novak()
        // Compartments and Species:
        compartment cytoplasm;
        species cyclin in cytoplasm, dimer in cytoplasm, dimer_p in cytoplasm, p_dimer in
      cytoplasm;
        species p_dimer_p in cytoplasm, cdc25_p in cytoplasm, wee1_p in cytoplasm;
        species IE_p in cytoplasm, UbE_star in cytoplasm, $cdc2 in cytoplasm, $cdc25 in cytoplasm;
        species $wee1 in cytoplasm, $IE in cytoplasm, $UbE in cytoplasm;
        // Assignment Rules:
        cdc2 := total_cdc2 - (dimer + p_dimer + p_dimer_p + dimer_p);
        cdc25 := total_cdc25 - cdc25_p;
        wee1 := total_wee1 - wee1_p;
        IE := total_IE - IE_p;
        UbE := total_UbE - UbE_star;
        k25 := V25_prime*(total_cdc25 - cdc25_p) + V25_double_prime*cdc25_p;
        kwee := Vwee_prime*wee1_p + Vwee_double_prime*(total_wee1 - wee1_p);
        k2 := V2_prime*(total_UbE - UbE_star) + V2_double_prime*UbE_star;
        total_cyclin := cyclin + dimer + dimer_p + p_dimer + p_dimer_p;
        Y15P := p_dimer + p_dimer_p;
        // Reactions:
        R1: => cyclin; k1AA;
        R2: cyclin => ; k2*cyclin;
        R3: cyclin + $cdc2 => dimer; k3*cyclin*cdc2;
        R4: dimer_p => dimer; kinh*dimer_p;
        R5: dimer => p_dimer; kwee*dimer;
        R6: dimer => dimer_p; kcak*dimer;
        R7: dimer => ; k2*dimer;
        R8: p_dimer => dimer; k25*p_dimer;
        R9: p_dimer => p_dimer_p; kcak*p_dimer;
        R10: p_dimer => ; k2*p_dimer;
        R11: p_dimer_p => p_dimer; kinh*p_dimer_p;
        R12: dimer_p => p_dimer_p; kwee*dimer_p;
        R13: p_dimer_p => dimer_p; k25*p_dimer_p;
        R14: p_dimer_p => ; k2*p_dimer_p;
        R15: dimer_p => ; k2*dimer_p;
        R17: $cdc25 => cdc25_p; ka*dimer_p*(total_cdc25 - cdc25_p)/(K_a + total_cdc25 - cdc25_p);
        R18: cdc25_p => $cdc25; kbPPase*cdc25_p/(K_b + cdc25_p);
        R19: $wee1 => wee1_p; ke*dimer_p*(total_wee1 - wee1_p)/(K_e + total_wee1 - wee1_p);
        R20: weel_p => $weel; kfPPase*weel_p/(K_f + weel_p);
        R21: $IE => IE_p; kg*dimer_p*(total_IE - IE_p)/(K_g + total_IE - IE_p);
        R22: IE_p => $IE; khPPAse*IE_p/(K_h + IE_p);
        R23: $UbE => UbE_star; kc*IE_p*(total_UbE - UbE_star)/(K_c + total_UbE - UbE_star);
        R24: UbE_star => $UbE; kd_anti_IE*UbE_star/(K_d + UbE_star);
        // Species initializations:
        cyclin = 100;
        dimer = 0;
        dimer_p = 0;
        p_dimer = 0;
        p_dimer_p = 0;
        cdc25_p = 0;
        wee1_p = 0;
        IE_p = 0;
        UbE_star = 0;
        // Compartment initializations:
        cytoplasm = 1;
        // Variable initializations:
        total_cdc2 = 100;
        total_cdc25 = 1;
        total_wee1 = 1;
        total_IE = 1;
        total_UbE = 1;
        V25_prime = 0.1;
        V25_double_prime = 2;
        Vwee_prime = 0.1;
        Vwee_double_prime = 1;
        V2_{prime} = 0.015;
        V2_double_prime = 1;
        k1AA = 1;
        k3 = 0.01;
        kinh = 0.025;
        kcak = 0.25;
        ka = 0.01;
        K_a = 0.1;
        kbPPase = 0.125;
        K_b = 0.1;
        ke = 0.0133;
        K_e = 0.3;
        kfPPase = 0.1;
        K_f = 0.3;
        kg = 0.0065;
        K_g = 0.01;
        khPPAse = 0.087;
        K_h = 0.01;
        kc = 0.1;
        K_c = 0.01;
        kd_anti_IE = 0.095;
        K_d = 0.01;
        // Other declarations:
        var k25, kwee, k2, total_cyclin, Y15P;
        const cytoplasm, total_cdc2, total_cdc25, total_wee1, total_IE, total_UbE;
        const V25_prime, V25_double_prime, Vwee_prime, Vwee_double_prime, V2_prime;
        const V2_double_prime, k1AA, k3, kinh, kcak, ka, K_a, kbPPase, K_b, ke;
        const K_e, kfPPase, K_f, kg, K_g, khPPAse, K_h, kc, K_c, kd_anti_IE, K_d;
        // Unit definitions:
        unit substance = 1e-9 mole;
        unit time_unit = 60 second;
        // Display Names:
        substance is "nanomole";
        time_unit is "minutes";
        dimer is "cyclin-cdc2 dimer";
        dimer_p is "Thr161 phosphorylated dimer(active MPF)";
        p_dimer is "Tyr15 phosphorylated dimer";
        p_dimer_p is "Thr161-Tyr15 phosphorylated dimer";
        cdc25_p is "phosphorylated cdc25";
        weel_p is "phosphorylated weel";
        IE_p is "phosphorylated intermediary enzyme";
        UbE_star is "ubiquitin conjugating enzyme";
        IE is "intermediary enzyme";
        UbE is "ubiquitin conjugating enzyme";
        R1 is "cyclin synthesis";
        R2 is "cyclin degradation";
        R3 is "cyclin-cdc2 dimer formation";
        R4 is "Thr161 dephosphorylation";
        R5 is "Tyr15 phosphorylation";
        R6 is "Thr161 phosphorylation";
        R7 is "cyclin degradation";
        R8 is "Tyr15 dephosphorylation";
        R9 is "Thr161 phosphorylation";
        R10 is "cyclin degradation";
        R11 is "Thr161 dephosphorylation";
        R12 is "Tyr15 phosphorylation";
        R13 is "Tyr15 dephosphorylation";
        R14 is "cyclin degradation";
        R15 is "cyclin degradation";
        R17 is "cdc25 activation";
        R18 is "cdc25 deactivation";
        R19 is "weel deactivation";
        R20 is "weel activation";
        R21 is "intermediary enzyme activation";
        R22 is "intermediary enzyme deactivation";
        R23 is "ubiquitin conjugating enzyme activation";
        R24 is "ubiquitin conjugating enzyme deactivation";
      end
      model_extract = model "novak"
      model_intact = model "novak" with k1AA = 1.8, V2_prime = 0.03, ka = 0.05, kc = 0.4, ke =
      0.0067, kg = 0.02, kbPPase = 0.0375, kd_anti_IE = 0.25, kfPPase = 0.05, khPPAse = 0.27
      sim1 = simulate uniform(0, 200, 1000)
      extract = run sim1 on model_extract
      intact = run sim1 on model_intact
      plot "M-phase Control, Extract (Novak, 1993)" extract.time vs extract.total_cyclin,
      extract.p_dimer_p, extract.dimer_p
      plot "M-phase Control, Intact Embryo (Novak, 1993)" intact.time vs intact.total_cyclin,
      intact.p_dimer_p, intact.dimer_p
                          M-phase Control, Extract (Novak, 1993)
             100
                                                                           extract.total_cyclin
                                                                           extract.p_dimer_p
                                                                           extract.dimer_p
              50
                                                                           intact.total_cyclin
                                                                           · intact.p_dimer_p
              0
                                                                         intact.dimer_p
                            50
                                         100
                                                      150
               0
                       M-phase Control, Intact Embryo (Novak, 1993)
```



100

50

0

50

100

150

200