

1 Overview

Note: This is the same code as in the quickstart guide. See the guide for a more in detail explanation.

```
1 from esbmtk import Model, Element, Species, Reservoir
2 from esbmtk import Signal, Connect, Source, Sink, Flux
3 from esbmtk import ExternalData
4
5 # create model
6 Model(
7     name="C_Cycle", # model name
8     stop=100,       # end time of model
9     time_unit="yr",  # time units
10    dt=1,            # time step
11 )
```

```
1 # Element properties
2 Element(
3     name="C",          # Element Name
4     model=C_Cycle,     # Model handle
5     mass_unit="mmol",  # base mass unit
6     li_label="C12",    # Name of light isotope
7     hi_label="C13",    # Name of heavy isotope
8     d_label="$\delta^{13}$C", # Name of isotope delta
9     d_scale="VPDB",     # Isotope scale. End of plot labels
10    r=0.0112372, # VPDB C13/C12 ratio https://www-pub.iaea.org/MTCD/publications/PDF/te\_825\_
11 )
12
13 # add species
14 Species(name="CO2", element=C) # Name & element handle
15 Species(name="DIC", element=C)
16 Species(name="OM", element=C)
17 Species(name="CaCO3", element=C)
```

```
1 Signal(name = "ACOR",
2         species = CO2,
3         duration = 100, # must match what is in the file
4         filename = "test-data.csv"
5 )
```

Once a signal instance has been created, it can be passed to a connector object in order to associate it with a flux (see the first connection below as an example).

```

1 Source(name="Fossil_Fuel_Burning", species=C02)
2 Source(name="Carbonate_Weathering", species=C02)
3 Source(name="Organic_Weathering", species=C02)
4 Source(name="Volcanic", species=C02)
5 Sink(name="Carbonate_burial", species=CaC03)
6 Sink(name="OM_burial", species=OM)
7
8 Reservoir(
9     name="Ocean",          # Name of reservoir
10    species=DIC,           # Species handle
11    delta=0,               # initial delta
12    concentration=2.62,    # concentration
13    unit="mmol",          # mass unit
14    volume=1.332E18,       # reservoir size (m^3)
15 )

```

```

1 # connect source to reservoir
2 Connect(
3     source=Fossil_Fuel_Burning, # source of flux
4     sink=Ocean,                 # target of flux
5     rate=0,                     # weathering flux in
6     delta=0,                    # set a default flux
7     pl=[ACOR],
8 )
9
10 Connect(
11     source=Carbonate_Weathering, # source of flux
12     sink=Ocean,                  # target of flux
13     rate=12.3E12,               # flux rate
14     delta=0,                    # isotope ratio
15 )
16
17 Connect(
18     source=Organic_Weathering, # source of flux
19     sink=Ocean,                 # target of flux
20     rate=4.0E12,                # flux rate
21     delta=-20,                  # isotope ratio
22 )

```

```

23 )
24
25 Connect(
26     source=Volcanic, # source of flux
27     sink=Ocean,      # target of flux
28     rate=6.0E12,     # flux rate
29     delta=-5,        # set a default isotope ratio
30 )
31
32 Connect(
33     source=Ocean,    # source of flux
34     sink=OM_burial,  # target of flux
35     rate=4.2E12,     # flux rate
36     alpha=-26.32,    # isotope fractionation
37 )
38
39 Connect(
40     source=Ocean,    # source of flux
41     sink=Carbonate_burial, # target of flux
42     rate=18.1E12,    # flux rate
43     delta=0,         # isotope fractionation
44 )

```

```

1 # Run the model
2 C_Cycle.run()
3
4 # plot the results
5 C_Cycle.plot_data()
6 C_Cycle.save_data()

```
