

# Lecture 11-13: Radioactive Decay

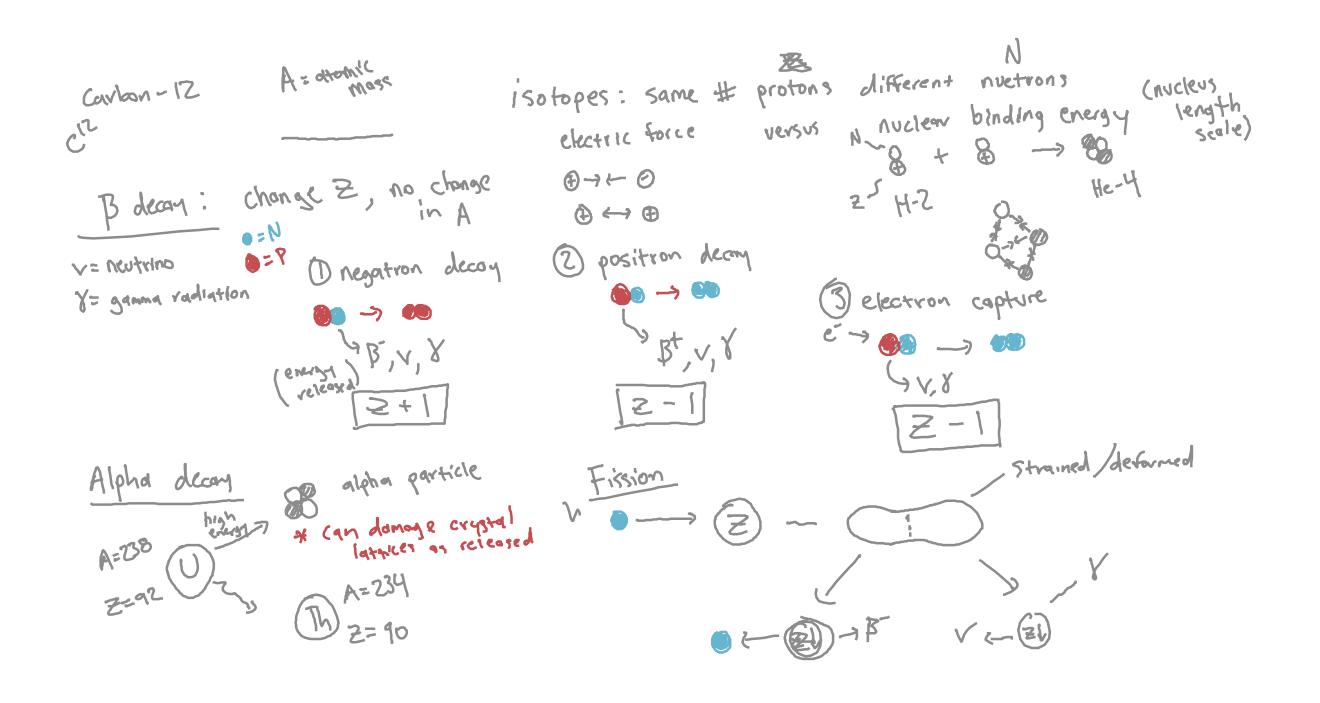
- 1. Mechanisms
- 2. The decay equation
  - A. Isochrons
  - B. Sm-Nd system
  - C. Pb-Pb dating (age of the Earth)

We acknowledge and respect the  $l \ni k^w \ni j \ni n$  peoples on whose traditional territory the university stands and the Songhees, Esquimalt and W S ANE E peoples whose historical relationships with the land continue to this day.





## Mechanisms of radioactive decay.



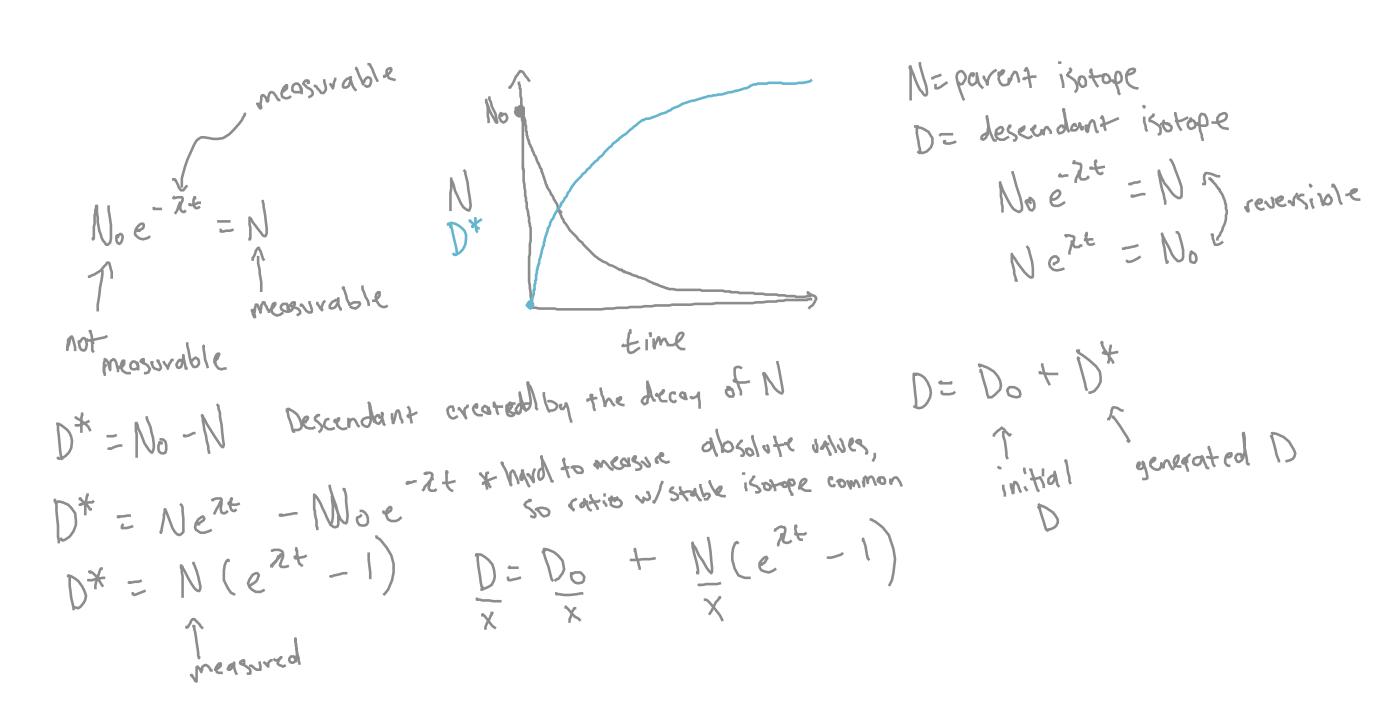




# The decay equation.

RutherFord + Soldy 1902 
$$N=$$
 number of moles of an isotope  $\frac{dN}{N}=\int -2dt$   $\ln x \ln dt$   $dx + 2dt$   $dx + 2dt$ 

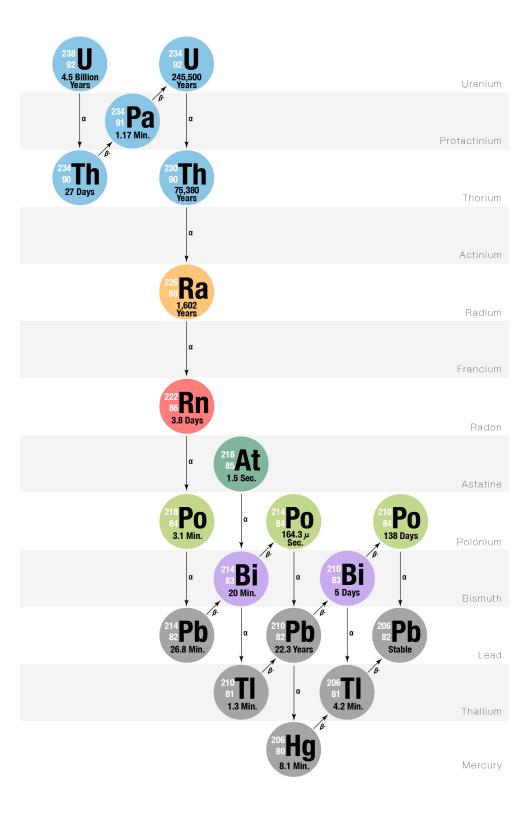
# The decay equation.





#### Half life

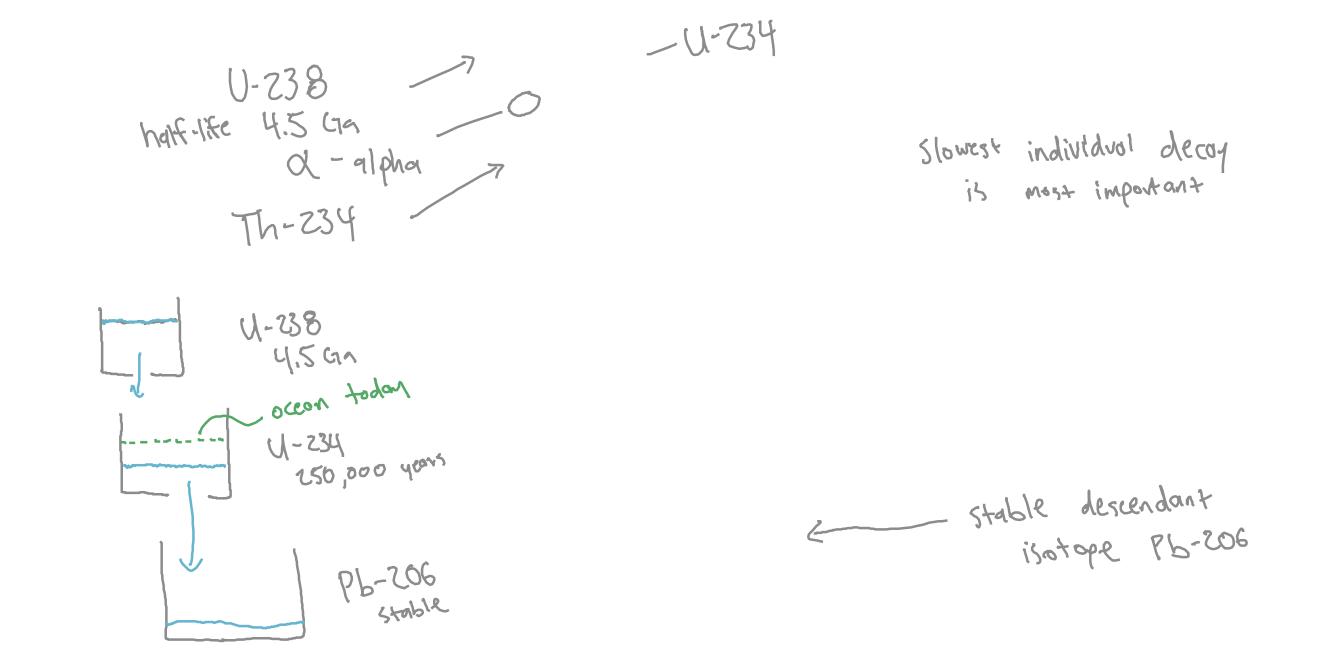


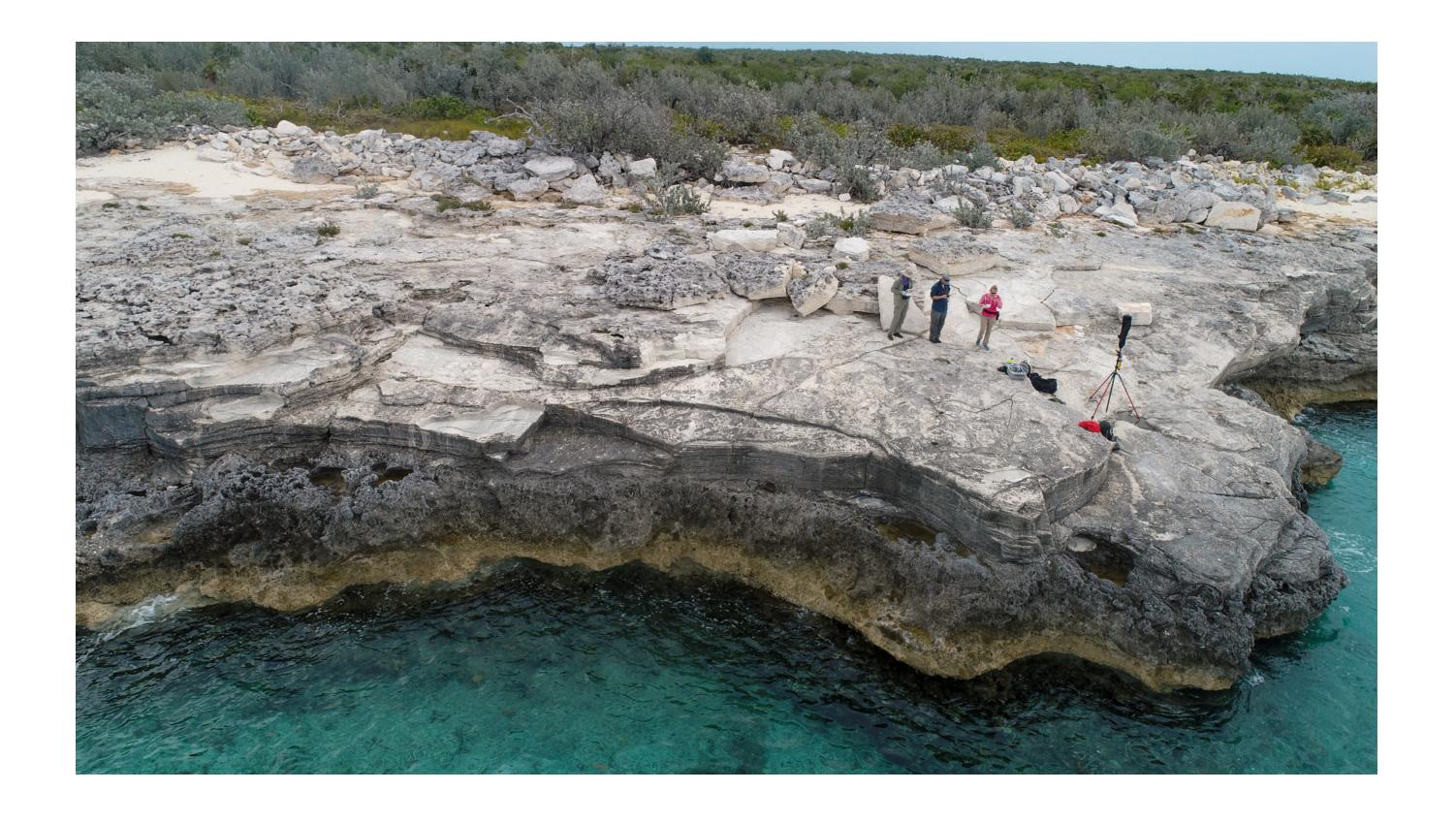






# U-Series dating of corals

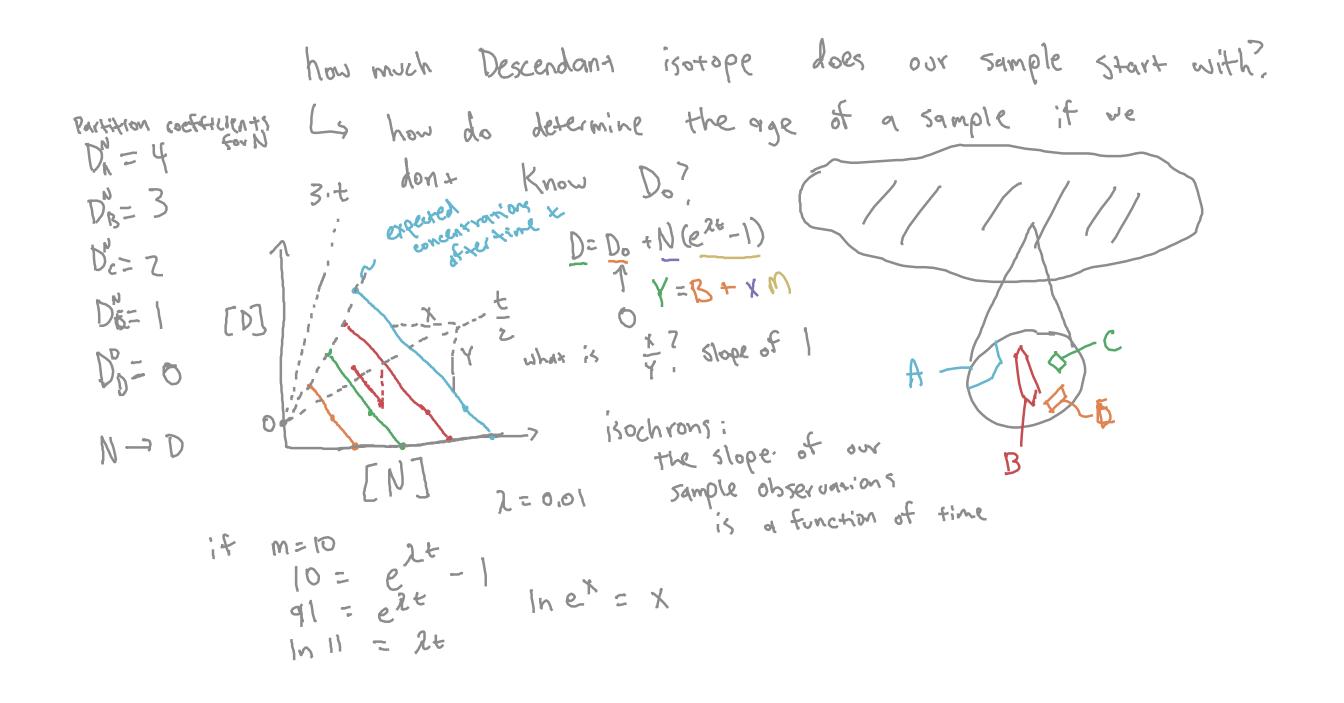








#### Isochrons





if a suite of samples forms a line in  $\frac{D}{N}$ , then the samples same age and initial Do. to use an isochron: 1. Samples have same age 2. Samples have the same initial Descendant Bottope 3. D and N isotopes to not enter of leave the mineral or rock since formation (closed 345+cm) it helps when the samples have very different N concentrations

Recall:

$$N_0 e^{-\lambda t} = N$$



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$$N_0 e^{-\lambda t} = N$$

```
In [3]: N_0 = 1 # parent isotope initial
         LAMBDA = 1e-4 # decay constant
         t = np.linspace(0, 1e5, 1000) #time (from now to 100 thousand years ago)
         N = N_0 * np.exp(-LAMBDA * t) # parent isotope
         D = N_0 - N #descendant isotope
```





#### Recall:

$$N_0 e^{-\lambda t} = N$$

```
In [3]: N 0 = 1 \# parent isotope initial
          LAMBDA = 1e-4 # decay constant
          t = np.linspace(0, 1e5, 1000) #time (from now to 100 thousand years ago)
          N = N_0 * np.exp(-LAMBDA * t) # parent isotope
          D = N_0 - N #descendant isotope
In [4]:
         plt.figure(figsize=(5, 5))
          plt.plot(t, N, label="Parent", alpha=1, lw=3)
          plt.plot(t, D, label="Descendant", alpha=1, lw=3)
          plt.legend(loc="best")
          _ = plt.gca().set_xlabel("Time")
          _ = plt.gca().set_ylabel("Amount")
                                                             1.0
                                                             8.0
                                                           9.0
0.4
                                                                              Parent
                                                                                Descendant
                                                             0.2
                                                             0.0
                                                                     25000 50000 75000 100000
                                                                           Time
```





Batch melting:

$$\frac{C_S}{C_0} = \frac{D}{F + D(1 - F)}$$



#### Batch melting:

$$\frac{C_S}{C_0} = \frac{D}{F + D(1 - F)}$$

```
In [10]:
    def batch_S(F, D, Co): #batch melting equation for solid
        return (D*Co) / (F + D * (1 - F))

D_n = np.array([1,2,3,4]) #a list of partition coefficients
F = 0.8 #melt fraction
N_0 = batch_S(F, D_n, 1) #C_1 for each F
```





#### Batch melting:

$$\frac{C_S}{C_0} = \frac{D}{F + D(1 - F)}$$

```
In [10]:
           def batch_S(F, D, Co): #batch melting equation for solid
               return (D*Co) / (F + D * (1 - F))
           D_n = np.array([1,2,3,4]) #a list of partition coefficients
           F = 0.8 \# melt fraction
           N_0 = batch_S(F, D_n, 1) \#C_1 for each F
In [11]:
           plt.figure(figsize=(6, 5))
           plt.plot(D_n, N_0, "o")
            _ = plt.gca().set_xlabel("Bulk partition coefficient D$_n$")
           _ = plt.gca().set_ylabel("N$_0$")
                                                               2.50
                                                               2.25
                                                               2.00
                                                             ≥° 1.75
                                                               1.50
                                                               1.25
                                                               1.00
                                                                             2
                                                                                       3
                                                                         Bulk partition coefficient D<sub>n</sub>
```





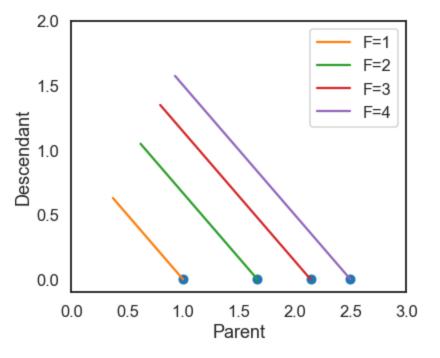
```
In [15]: N = [n * np.exp(-LAMBDA * t[:100])  for n in N_0] # the decay equation for each starting N
           C = [(n - NS) \text{ for } n, NS \text{ in } zip(N_0, N)] # the descendant isotope (initial minus current)
```





```
In [15]: N = [n * np.exp(-LAMBDA * t[:100]) for n in N_0] # the decay equation for each starting N C = [(n - NS)] for n, NS in zip(N_0, N) # the descendant isotope (initial minus current)
```

```
In [17]: plt.figure(figsize=(6, 5))
    plt.plot(N_0, [0, 0, 0, 0], "o", alpha=1)
    _ = [plt.plot(a, b, label="F=" + str(c), alpha=1) for a, b, c in zip(N, C, D_n)]
    plt.legend(loc="best")
    _ = plt.gca().set_xlabel("Parent")
    _ = plt.gca().set_ylabel("Descendant")
    _ = plt.gca().set_xlim([0, 3])
    _ = plt.gca().set_ylim([-0.1, 2])
```







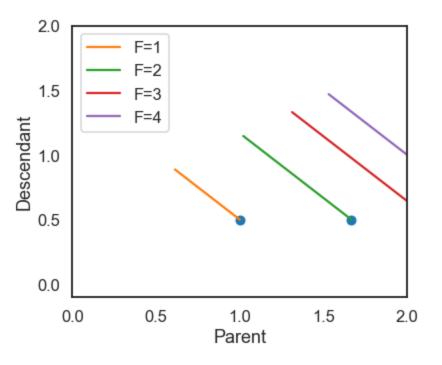
```
In [18]: | np.random.seed(12)
          N = [n * np.exp(-LAMBDA * t[:50]) for n in N_0] # the decay equation for each starting N
          Ic = 0.5*np.ones(4) # some fixed amount of initial child isotope
          C = [i + (n - NS)  for i, n, NS in zip(Ic, N_0, N)] # the descendant isotope (initial parent minus current)
```





```
In [18]:
    np.random.seed(12)
    N = [n * np.exp(-LAMBDA * t[:50]) for n in N_0] # the decay equation for each starting N
    Ic = 0.5*np.ones(4) # some fixed amount of initial child isotope
    C = [i + (n - NS) for i, n, NS in zip(Ic, N_0, N)] # the descendant isotope (initial parent minus current)
```

```
In [19]: plt.figure(figsize=(6, 5))
    plt.plot(N_0, Ic, "o")
    _ = [plt.plot(a, b, label="F=" + str(c)) for a, b, c in zip(N, C, D_n)]
    plt.legend(loc="best")
    _ = plt.gca().set_xlabel("Parent")
    _ = plt.gca().set_ylabel("Descendant")
    _ = plt.gca().set_xlim([0, 2])
    _ = plt.gca().set_ylim([-0.1, 2])
```



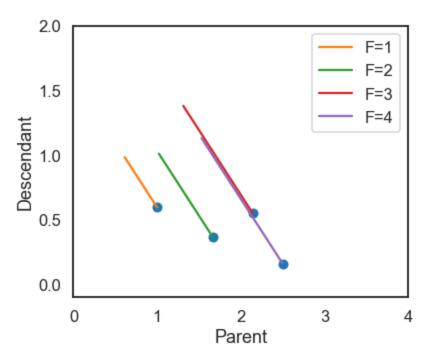






```
In [20]: Ic = 0.5 + np.array(np.random.normal(0.0, 0.2, 4)) # some random amount of initial child isotope
          C = [i + (n - NS)  for i, n, NS in zip(Ic, N_0, N)] # the descendant isotope (initial parent minus current)
```

```
In [22]: plt.figure(figsize=(6, 5))
          plt.plot(N_0, Ic, "o")
          _ = [plt.plot(a, b, label="F=" + str(c)) for a, b, c in zip(N, C, D_n)]
          plt.legend(loc="best")
           _ = plt.gca().set_xlabel("Parent")
           _ = plt.gca().set_ylabel("Descendant")
           _ = plt.gca().set_xlim([0, 4])
          _ = plt.gca().set_ylim([-0.1, 2])
```







### **Utility**

Review: any of our melting models

could produce variations

in Pisotopes

isochrons:  $V=M\times +b$  — the deay equation reduces to this form

with the following assumptions:

M= ext2 -1

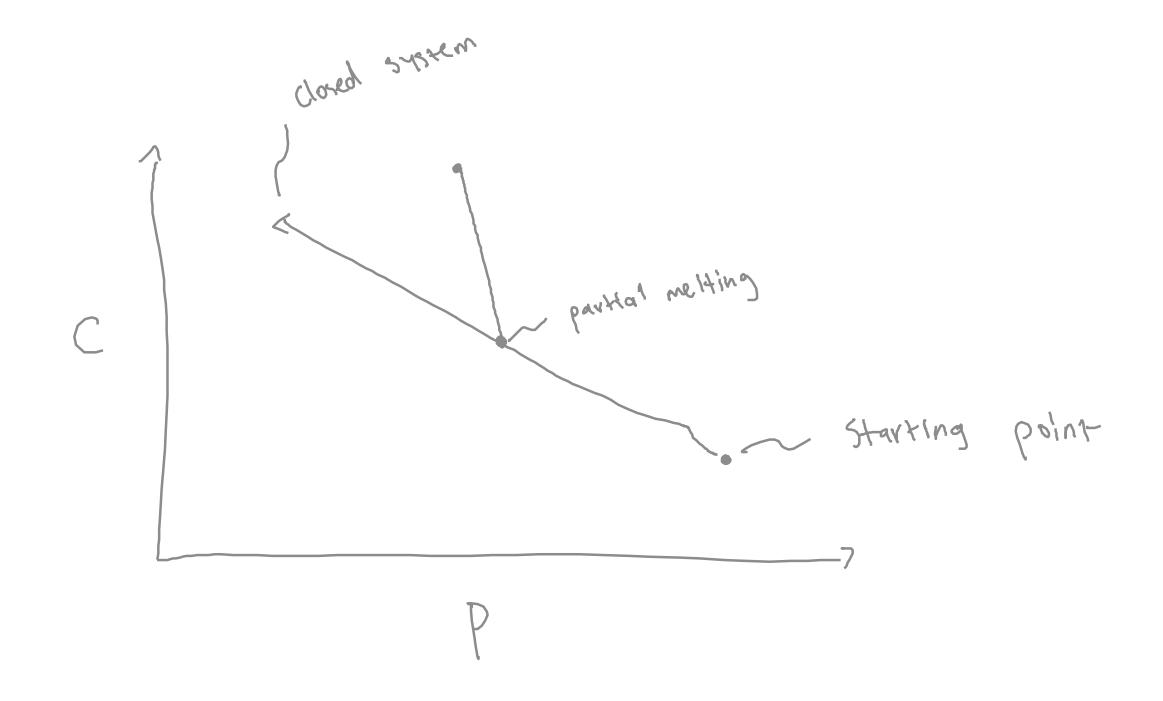
Same initial Daughter isotope concentration

- · same age of samples

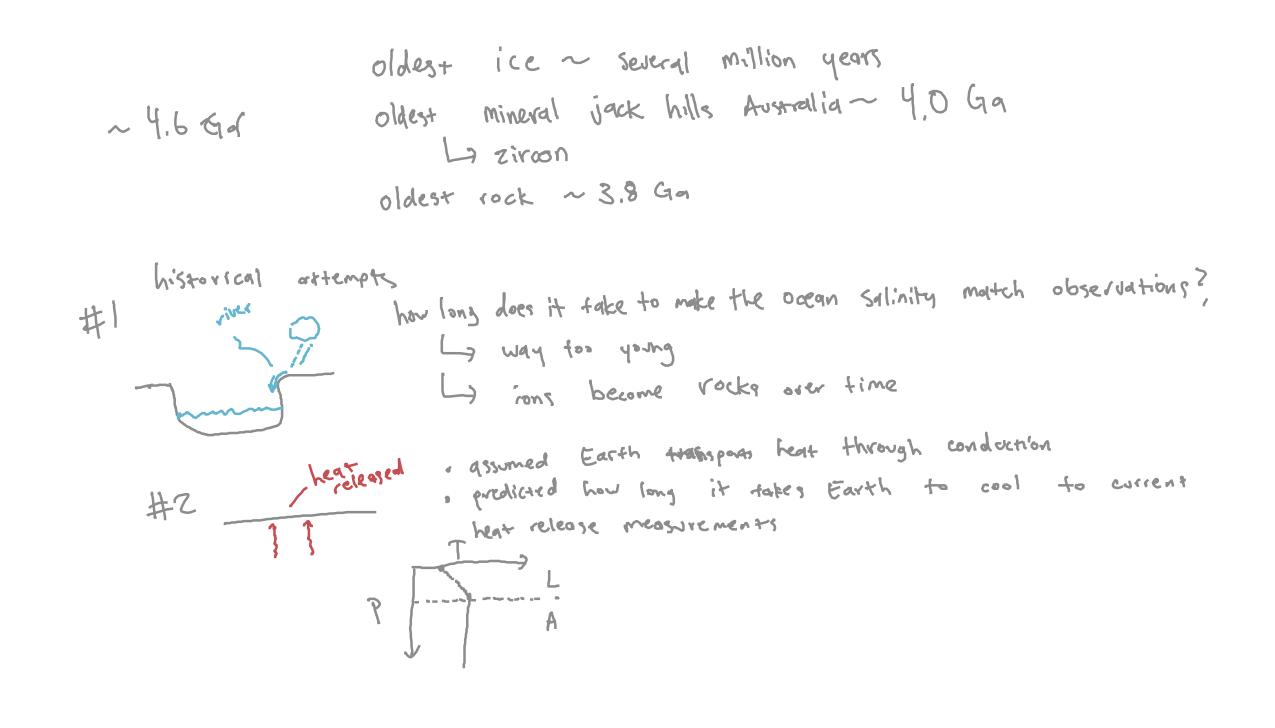




Sm-Nd decay and model ages (the chondritic uniform reservoir or CHUR)  $^{147}$  Sm decays to  $^{143}$  Nd through alpha decay with a decay constant of  $\lambda^{147}$  =  $6.54\times10^{-12}$ 

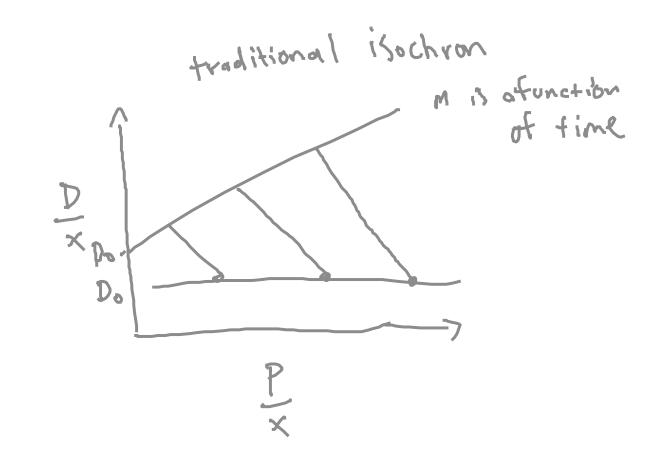


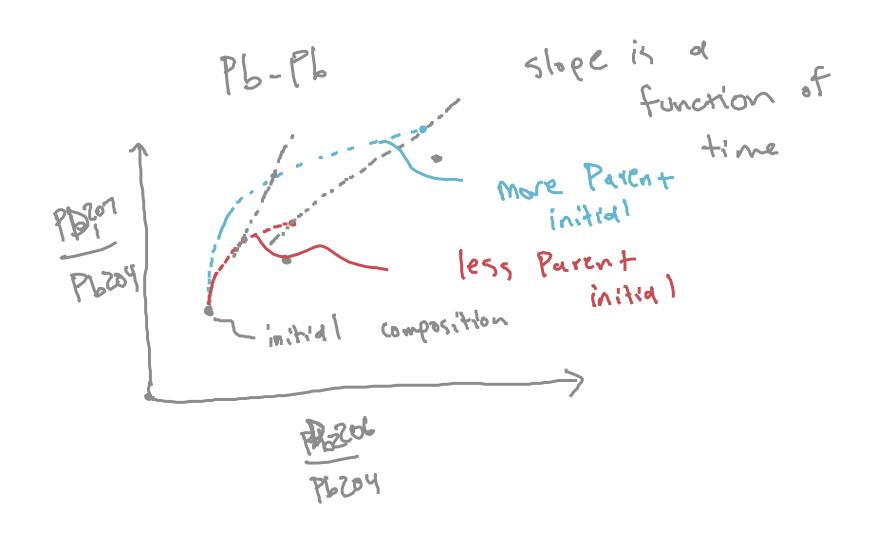














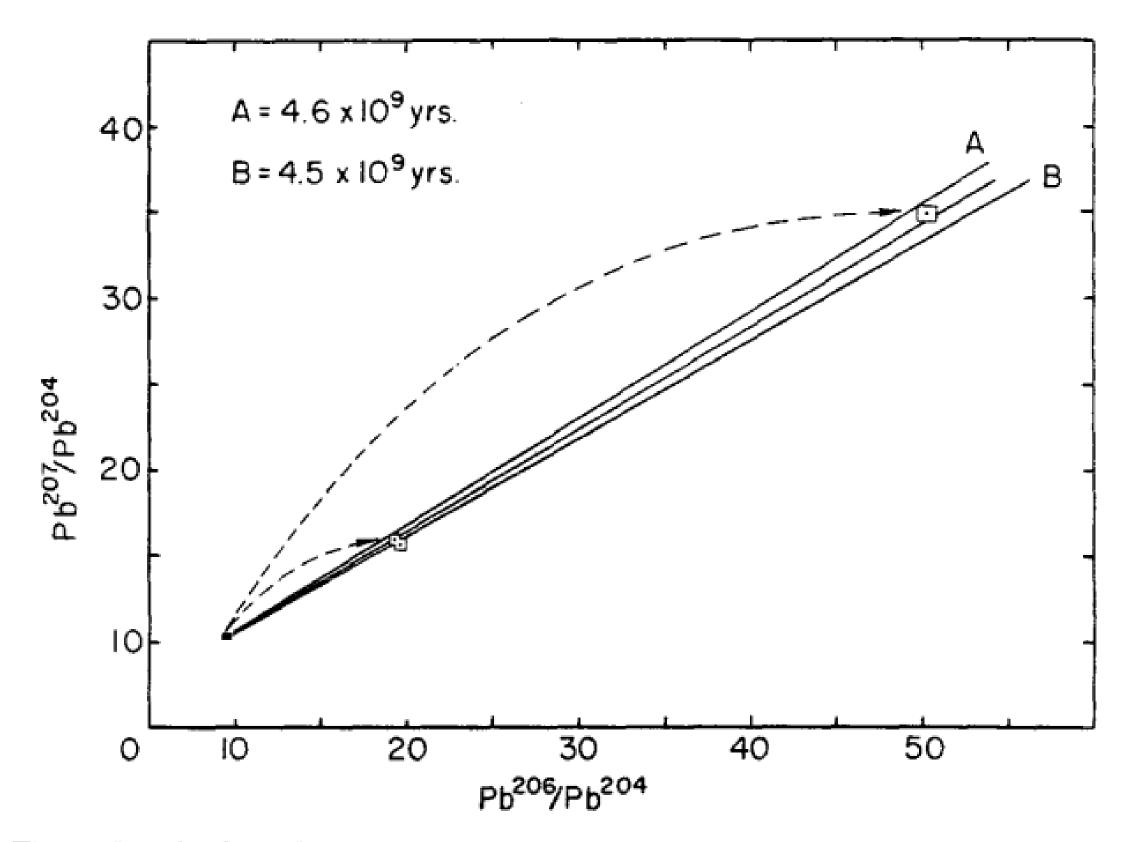


Fig. 1. The lead isochron for meteorites and its estimated limits. The outline around each point indicates measurement error.



