

Paschen Curve Monte Carlo Simulations

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Basic 1D parallel plate simulations

```
In[=]:= AvgElectronsBanked[Nc_, Ni_, count_] := Module[{result, stack},
  stack = CreateDataStructure["Stack"];
  result = Table[
    Module[{pos1, pos2, numElectrons, lambdaI},
      numElectrons = 1;
      stack["Push", 0];
      While[! stack["EmptyQ"],
        pos1 = stack["Pop"];
        While[(pos2 = pos1 + RandomVariate[ExponentialDistribution[1]]) ≤ Nc,
          lambdaI = Nc / Ni;
          If[(pos2 - pos1) ≥ lambdaI, numElectrons++];
          stack["Push", pos2]];
        Sow[pos2];
        pos1 = pos2];
      numElectrons],
    {count}];
  Around[Mean[result], StandardDeviation[result] / Sqrt[count]]]
```

```
In[5]:= AvgElectronsWeighted[Nc_, Ni_, count_] := Module[{result},
  result = Table[
    Module[{pos1, pos2, numElectrons, lambdaI, weight},
      numElectrons = 1;
      pos1 = 0;
      weight = 1;
      While[(pos2 = pos1 + RandomVariate[ExponentialDistribution[1]]) ≤ Nc,
        lambdaI = Nc / Ni;
        If[(pos2 - pos1) ≥ lambdaI, numElectrons += weight;
          weight *= 2];
        pos1 = pos2];
      numElectrons],
    {count}];
  Around[Mean[result], StandardDeviation[result] / Sqrt[count]]]
```

Defining 3D MC utility functions

```
In[6]:= GetTime[s_, vx_, vw_, ax_] := Module[{S, set, start},
  If[vw > 0,
    S[t_] := 1 / (2 ax) * ((vx + ax t) Sqrt[vw^2 + (vx + ax t)^2] -
      vx Sqrt[vw^2 + vx^2] + vw^2 (ArcSinh[(ax t + vx) / vw] - ArcSinh[vx / vw])),
    S[t_] := (-vx Sqrt[vx^2] + (ax t + vx) Sqrt[(ax t + vx)^2]) / (2 ax)
  ];
  set = FindRoot[S[t] == s, {t, 0.05, 10}];
  t /. set]

In[7]:= GetVelocity[energy_, cosθ_, φ_] := Module[{q, m, eθ, v, vx1, vy1, vz1},
  q = 1;
  m = 511000;
  v = Sqrt[2 energy / m];
  vx1 = v Cos[φ] Sqrt[1 - cosθ^2];
  vy1 = v Sin[φ] Sqrt[1 - cosθ^2];
  vz1 = v cosθ;
  {vx1, vy1, vz1}]
```

```

In[6]:= GetNewPositions[E_, s_, x0_, y0_, z0_, vx_, vy_, vz_] :=
  Module[{vw, q, m, ax, set, x, y, z, t, velocities},
    vw = Sqrt[vy^2 + vz^2];
    q = 1;
    m = 511000;
    ax = (q E) / m;
    t = GetTime[s, vx, vw, ax];
    x = vx t + 1/2 ax t^2;
    y = vy t;
    z = vz t;
    {x0 + x, y0 + y, z0 + z, vx + ax t, vy, vz, 1/2 * m * ((vx + ax t)^2 + vy^2 + vz^2)}
  ]

In[7]:= UpdateVelocity[energy_, vx_, vy_, vz_, tbound_] :=
  Module[{m, cost, phi, a, pvec1, pvec2, pvec3, b, c, v, res},
    If[energy == 0, Return[{0, 0, 0, 100, 100}]];
    m = 511000;
    cost = RandomReal[{tbound, 1}, WorkingPrecision → 16];
    phi = RandomReal[2 Pi, WorkingPrecision → 16];
    a = Normalize[{vx, vy, vz}];
    pvec1 = Normalize[{-vy, -vx, -0}];
    pvec2 = Normalize[{0, -vz, vy}];
    pvec3 = Normalize[{-vz, 0, vx}];
    b = pvec1;
    If[Abs[Dot[a, pvec2]] < Abs[Dot[a, b]], b = pvec2];
    If[Abs[Dot[a, pvec3]] < Abs[Dot[a, b]], b = pvec3];
    c = Cross[a, b];
    v = Sqrt[2 energy / m];
    res = SetPrecision[
      v (a cost + b Sqrt[1 - cost^2] Cos[phi] + c Sin[phi] Sqrt[1 - cost^2]), 16];
    Flatten[Append[res, {cost, phi}]]
  ]

```

```

In[5]:= GetNewPositionsEuler[Δt_, s_, V_, x0_, y0_, z0_, vx0_, vy0_, vz0_] :=
Module[{q, m, path, eField, R, x, y, z, vx, vy,
vz, ex, ey, ez, e0, e1, Δpos, Δr, Δtsum, scalefactor},
q = 1;
m = 511000;
{x, y, z} = {x0, y0, z0};
{vx, vy, vz} = {vx0, vy0, vz0};
path = 0;
Δtsum = 0;
While[path < s,
(*check energy conservation with delta t - take rms error *)
(*Sow[{Δtsum,x,y,z,vx,vy,vz,ez,ey,ez}];*)
Δtsum += Δt;
Δpos = Δt {vx, vy, vz};
Δr = Norm[Δpos];
If[path + Δr > s ,
scalefactor = (s - path) / Δr;
Δpos *= scalefactor;
{x, y, z} += Δpos;
path += Norm[Δpos];
If[! InBounds[x, y, z], Return[{x, y, z, vx, vy, vz, 0, 0, 0, 0}]];
{ex, ey, ez} = GetEField[V, x, y, z];
{vx, vy, vz} += ((q (Δt * scalefactor)) / m) {ex, ey, ez};
Break[]];
];
{x, y, z} += Δpos;
path += Norm[Δpos];
If[! InBounds[x, y, z], Return[{x, y, z, vx, vy, vz, 0, 0, 0, 0}]];
{ex, ey, ez} = GetEField[V, x, y, z];
{vx, vy, vz} += ((q Δt) / m) {ex, ey, ez};
];
(*Sow[{Δtsum,x,y,z,vx,vy,vz}];*)
e0 = 1/2 * m * (vx0^2 + vy0^2 + vz0^2) +
GetPotential[V, x0, y0, z0] - GetPotential[V, x, y, z];
e1 = 1/2 * m * (vx^2 + vy^2 + vz^2);
{x, y, z, vx, vy, vz, e1, e1 - e0, path, s}]

```

Defining utility functions for electrode geometries

```

In[=]:= (*Parallel plate*)
d = 5;
ID = "PP";
GetStartingPos[] := {0, 0, 0, 0, 0, 0};
InBounds[x_, y_, z_] := 0 ≤ x ≤ d;
GetEField[V_, x_, y_, z_] := {V / d, 0, 0};
GetPotential[V_, x_, y_, z_] := V x / d;

In[=]:= (*Concentric spheres*)
a = 2;
b = 7;
ID = "SS";
GetStartingPos[] := Flatten[
  Append[RandomPoint[Sphere[{0, 0, 0}, a], 1, WorkingPrecision → 16], {0, 0, 0}]];
InBounds[x_, y_, z_] := a ≤ Norm[{x, y, z}] ≤ b
(*k/r^2 → k = Vab/(b-a)*)
GetEField[V_, x_, y_, z_] :=
  {x, y, z} (V a b) / (Norm[SetPrecision[{x, y, z} // N, 16]]^3 (b - a));
(*GetEField[V_,x_,y_,z_] := {x,y,z} (V a b)/(Norm[{x,y,z}]^3 (b-a)); *)
GetPotential[V_, x_, y_, z_] := (V a b) / (Norm[{x, y, z}] (b - a));

```

```

ln[6]:= (*Concentric spheres discrete*)
sspotential = Import["/Users/wbowers/Downloads/phi2.txt", "Table"];
sspotential = Drop[sspotential, 2];
a = 2;
b = 7;
ID = "SS Discrete";
GetStartingPos[] := Flatten[
  Append[RandomPoint[Sphere[{0, 0, 0}, a], 1, WorkingPrecision → 16], {0, 0, 0}]];
InBounds[x_, y_, z_] := a ≤ Norm[{x, y, z}] ≤ b
gs = 0.01;
GetPotential[V_, x_, y_, z_] := Module[{ρ, length, i, j, α, β},
  ρ = Sqrt[x^2 + y^2];
  i = Floor[ρ / gs];
  j = Floor[Abs[z] / gs];
  length = Length[sspotential];
  If[i + 2 > length || j + 2 > length, Return[0.]];
  α = ρ / gs - i;
  β = Abs[z] / gs - j;
  -V ((1 - β) (1 - α) sspotential[[i + 1, j + 1]] + (1 - β) α sspotential[[i + 2, j + 1]] +
    β (1 - α) sspotential[[i + 1, j + 2]] + β α sspotential[[i + 2, j + 2]])
]
GetEField[V_, x_, y_, zraw_] :=
Module[{z, ρ, i, j, length, α, β, Er, Ez, Eρ, Ezf, phi, flip},
  ρ = Sqrt[x^2 + y^2];
  i = Floor[ρ / gs];
  If[zraw < 0, z = Abs[zraw]; flip = True, z = zraw];
  j = Floor[z / gs];
  length = Length[sspotential];
  If[i + 2 > length || j + 2 > length, Return[{0., 0., 0.}]];
  α = ρ / gs - i;
  β = z / gs - j;
  Eρ = (1 - β) (sspotential[[i + 1, j + 1]] - sspotential[[i + 2, j + 1]]) / gs +
    β (sspotential[[i + 1, j + 2]] - sspotential[[i + 2, j + 2]]) / gs;
  Ezf = (1 - α) (sspotential[[i + 1, j + 1]] - sspotential[[i + 1, j + 2]]) / gs +
    α (sspotential[[i + 2, j + 1]] - sspotential[[i + 2, j + 2]]) / gs;
  If[flip, Ezf = -Ezf];
  -V Flatten[{Eρ * Normalize[{x, y}], Ezf}]
]

```

```

GetPotentialContinuous[V_, x_, y_, z_] := Module[{ρ, length, i, j, α, β},
  ρ = Sqrt[x^2 + y^2];
  i = Floor[ρ / gs];
  j = Floor[Abs[z] / gs];
  length = Length[sspotential];
  If[i + 2 > length || j + 2 > length, Return[0.]];
  α = ρ / gs - i;
  β = Abs[z] / gs - j;
  -V ((1 - β) (1 - α) sspotential[[i + 1, j + 1]] + (1 - β) α sspotential[[i + 2, j + 1]] +
    β (1 - α) sspotential[[i + 1, j + 2]] + β α sspotential[[i + 2, j + 2]])
]

GetEFieldContinuous[V_, x_, y_, zraw_] :=
Module[{z, ρ, i, j, length, α, β, Er, Ez, Eρ, Ezf, phi, flip},
  ρ = Sqrt[x^2 + y^2];
  i = Floor[ρ / gs];
  If[zraw < 0, z = Abs[zraw]; flip = True, z = zraw];
  j = Floor[z / gs];
  length = Length[sspotential];
  If[i + 2 > length || j + 2 > length, Return[{0., 0., 0.}]];
  α = ρ / gs - i;
  β = z / gs - j;
  Eρ = (1 - β) (sspotential[[i + 1, j + 1]] - sspotential[[i + 2, j + 1]]) / gs +
    β (sspotential[[i + 1, j + 2]] - sspotential[[i + 2, j + 2]]) / gs;
  Ezf = (1 - α) (sspotential[[i + 1, j + 1]] - sspotential[[i + 1, j + 2]]) / gs +
    α (sspotential[[i + 2, j + 1]] - sspotential[[i + 2, j + 2]]) / gs;
  If[flip, Ezf = -Ezf];
  -V Flatten[{Eρ * Normalize[{x, y}], Ezf}]
]

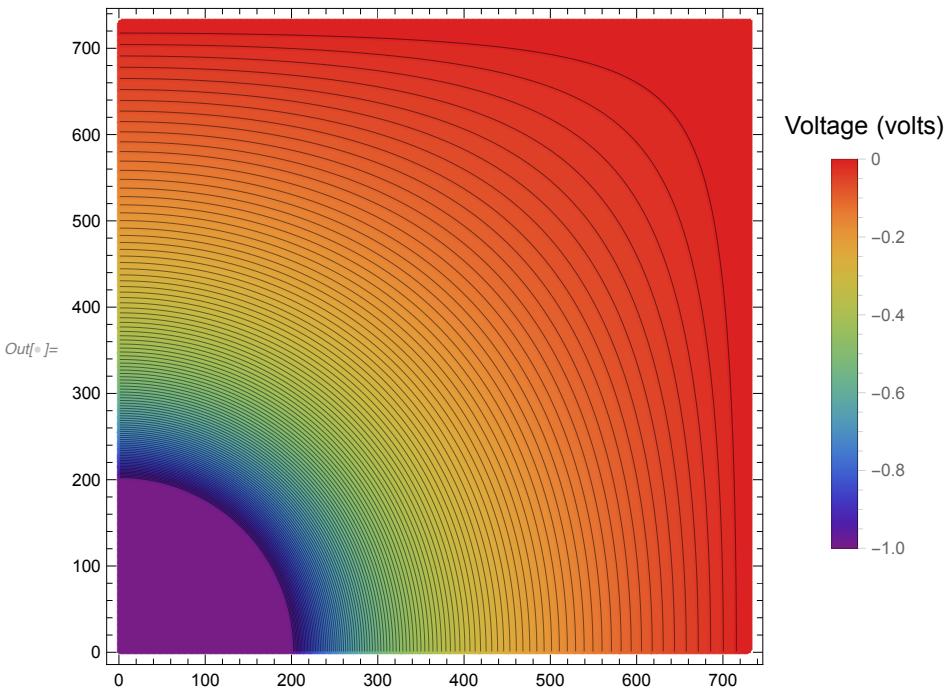
```

```

ln[6]:= (*Sphere in cylinder*)
scpotential = Import["/Users/wbowers/Downloads/new-cwPhi.txt", "Table"];
GetScPotential[i_, j_] :=
  If[i > Length[scpotential] || j > Length[scpotential], 0, scpotential[[i, j]]];
a = 2;
b = 7.3;
h = 14.6;
ID = "SC";
lmax = 10;
binwidth = 0.05;
gs = 100 * scpotential[[3, 3]];
phiFile = "Phi file: " <> StringDelete[ToString[scpotential[[1]]], {"", {"", ""}}];
scpotential = Drop[scpotential, 3];
GetAngles[x_, y_, z_] := {z / Sqrt[x^2 + y^2 + z^2], ArcTan[x, y]};
FindBinNum[costheta_, binwidth_] := Floor[(costheta + 1) / binwidth] + 1;
GetStartingPos[] := Flatten[
  Append[RandomPoint[Sphere[{0, 0, 0}, a], 1, WorkingPrecision \rightarrow 16], {0, 0, 0}]];
InBounds[x_, y_, z_] :=
  a \leq Sqrt[x^2 + y^2 + z^2] && Sqrt[x^2 + y^2] \leq b && -h / 2 \leq z \leq h / 2;
GetPotential[V_, x_, y_, z_] := Module[{rho, length, i, j, alpha, beta},
  rho = Sqrt[x^2 + y^2];
  i = Floor[Abs[z] / gs];
  j = Floor[rho / gs];
  alpha = Abs[z] / gs - i;
  beta = rho / gs - j;
  -V ((1 - beta) (1 - alpha) GetScPotential[i + 1, j + 1] + (1 - beta) alpha GetScPotential[i + 2, j + 1] +
    beta (1 - alpha) GetScPotential[i + 1, j + 2] + beta alpha GetScPotential[i + 2, j + 2])
]
GetEField[V_, x_, y_, zraw_] :=
  Module[{z, rho, i, j, length, alpha, beta, Er, Ez, Erho, Ezf, phi, flip},
  rho = Sqrt[x^2 + y^2];
  If[zraw < 0, z = Abs[zraw]; flip = True, z = zraw];
  i = Floor[z / gs];
  j = Floor[rho / gs];
  alpha = z / gs - i;
  beta = rho / gs - j;
  Erho = (1 - alpha) (GetScPotential[i + 1, j + 1] - GetScPotential[i + 1, j + 2]) / gs +
    alpha (GetScPotential[i + 2, j + 1] - GetScPotential[i + 2, j + 2]) / gs;
  Ezf = (1 - beta) (GetScPotential[i + 1, j + 1] - GetScPotential[i + 2, j + 1]) / gs +
    beta (GetScPotential[i + 1, j + 2] - GetScPotential[i + 2, j + 2]) / gs;
  If[flip, Ezf = -Ezf];
  -V Flatten[{Erho * Normalize[{x, y}], Ezf}]
]

```

```
In[5]:= fig5 = ListContourPlot[scpotential, Contours -> 100,
PlotLegends -> Placed[BarLegend[{"Rainbow", {-1, 0}}],
LegendLabel -> "Voltage (volts)", Automatic], ColorFunction -> "Rainbow"]
```



```
In[6]:= Export["/Users/wbowers/Documents/stem-fellowship/tif-figs/e-field-contour.TIFF",
fig5, ImageResolution -> 600]
```

```
Out[6]= /Users/wbowers/Documents/stem-fellowship/tif-figs/e-field-contour.TIFF
```

Creating e-field export for sphere-in-sphere

```
In[7]:= ssregion = RegionDifference[Ball[{0, 0, 0}, 7], Ball[{0, 0, 0}, 2]];
```

```
In[8]:= ssregion = RegionDifference[Ball[{0, 0, 0}, 7], Ball[{0, 0, 0}, 2]];
```

```
Timing[spherepotential =
```

```
NDSolveValue[{lapl, DirichletCondition[f[x, y, z] == -1, x^2 + y^2 + z^2 < 9],
```

```
DirichletCondition[f[x, y, z] == 0, x^2 + y^2 + z^2 ≥ 9]}, f, {x, y, z} ∈ ssregion]]
```

```
sspotential = Table[Module[{val},
```

```
val = SetPrecision[spherepotential[r, 0, z], 4];
```

```
If[val === Indeterminate && Sqrt[r^2 + z^2] ≤ 2, val = -1];
```

```
If[val === Indeterminate && Sqrt[r^2 + z^2] ≥ 6, val = 0];
```

```
SetPrecision[val, 16]
```

```
], {z, 0, 7, 0.008}, {r, 0, 7, 0.008}];
```

```
In[9]:= sspotential = Insert[sspotential,
```

```
"12/20/2022, William Bowers, a 1/r sphere-in-sphere potential", 1];
```

```
sspotential = Insert[sspotential, "701, 701, 0.0001", 2];
```

```
In[5]:= Export["/Users/wbowers/Downloads/phi2.txt", SetPrecision[sspotential, 4], "Table"]
Out[5]= /Users/wbowers/Downloads/phi2.txt
```

The main MC simulation (for parallel plate and sphere-in-sphere)

```
In[6]:= AvgElectrons3D[Nc_, Ni_, D_, Ui_, tbound_, count_] :=
Module[{V, E, λ, λi, result, stack},
λ = D / Nc;
λi = D / Ni;
V = Ui * Ni;
E = V / D;
stack = CreateDataStructure["Stack"];
result = Table[
Module[{x0, x1, y0, y1, z0, z1, numElectrons, cosθ, φ, s, vx, vy, vz, energy},
numElectrons = 1;
stack["Push", {0, 0, 0, 0, 0, 0}];
While[! stack["EmptyQ"],
{x0, y0, z0, vx, vy, vz} = stack["Pop"];
cosθ = 1;
φ = 0;
While[
s = RandomVariate[ExponentialDistribution[1 / λ]];
{x1, y1, z1, vx, vy, vz} =
GetNewPositions[E, s, x0, y0, z0, vx, vy, vz];
0 ≤ x1 ≤ D,
If[energy ≥ Ui,
numElectrons++;
stack["Push", {x1, y1, z1, 0, 0, 0}];
energy -= Ui,
(*else*)
energy -= Ui * RandomVariate[UniformDistribution[{0.1, 0.5}]];
If[energy < 0, energy = 0];
];
{vx, vy, vz} = UpdateVelocity[energy, vx, vy, vz, tbound];
{x0, y0, z0} = {x1, y1, z1};
];
];
];
numElectrons],
{count}];

Around[Mean[result], StandardDeviation[result] / Sqrt[count]]]

In[7]:= AvgElectronsEuler[Nc_, Ni_, D_, Ui_, dt_, tbound_, print_, count_] :=
Module[{eErrorList, V, E, λ, λi, result, stack, collisions, paths, svals, countbin,
```

```

electronsbin, squaredbin, lbin, lbinSquared, data, seedcount, totalpaths},
 $\lambda = D / N_c;$ 
 $\lambda_i = D / N_i;$ 
 $V = U_i * N_i;$ 
totalpaths = 0;
paths = {};
svals = 0;
E = V / D;
seedcount = 1;
collisions = {};
eErrorList = {};
stack = CreateDataStructure["Stack"];
result = Table[
  Module[{x0, x1, y0, y1, z0, z1, numElectrons, cosθ, φ,
    s, vx, vy, vz, energy, eError, numCollisions, pathval, sval, idx,
    cθ0, φ0, SH, l, l2, collisionE, cost, phi, start, path, input},
    numElectrons = 1;
    numCollisions = 0;
    stack["Push", GetStartingPos[]];
    {cθ0, φ0} = GetAngles[x0, y0, z0];
    While[! stack["EmptyQ"],
      {x0, y0, z0, vx, vy, vz} = stack["Pop"];
      cosθ = 1;
      φ = 0;
      energy = 0;
      path = 0;
      While[
        s = RandomVariate[ExponentialDistribution[1 / λ], WorkingPrecision → 16];
        {x1, y1, z1, vx, vy, vz, energy, eError, pathval, sval} =
          GetNewPositionsEuler[dt, s, V, x0, y0, z0, vx, vy, vz];
        AppendTo[eErrorList, eError];
        InBounds[x1, y1, z1],
        path += pathval;
        AppendTo[paths, pathval];
        svals += s;
        numCollisions++;
        If[energy ≥ U_i,
          numElectrons++;
          stack["Push", {x1, y1, z1, 0, 0, 0}];
          energy -= U_i,
          collisionE =
            SetPrecision[U_i * RandomVariate[UniformDistribution[{1 / 10, 1 / 2}], 16];
          energy -= collisionE;
          If[energy < 0, energy = 0];
        ];
      ];
    ];
  ];
];

```

```

];
{vx, vy, vz, cost, phi} = UpdateVelocity[energy, vx, vy, vz, tbound];
{x0, y0, z0} = {x1, y1, z1};
];
totalpaths += path;
];
AppendTo[collisions, numCollisions];
numElectrons,
{count}];
If[print,
Print["1000 * energy error error: ", 1000 * RootMeanSquare[eErrorList]];
Print["V/NC: ", V / Nc];
Print["λ: ", λ];
Print["Geometry: ", ID];
Print["Collisions: ",
Around[Mean[collisions], StandardDeviation[collisions] / Sqrt[count]]];
Print["Mean path: ", totalpaths / count / λ];
];
Around[Mean[result], StandardDeviation[result] / Sqrt[count]]]

```

Exploring discrete sphere-in-cylinder E field

```

In[=]:= lpc = D[ϕ[r, z], z, z] + D[ϕ[r, z], r, r] + 1/r D[ϕ[r, z], r];
TraditionalForm[lpc]

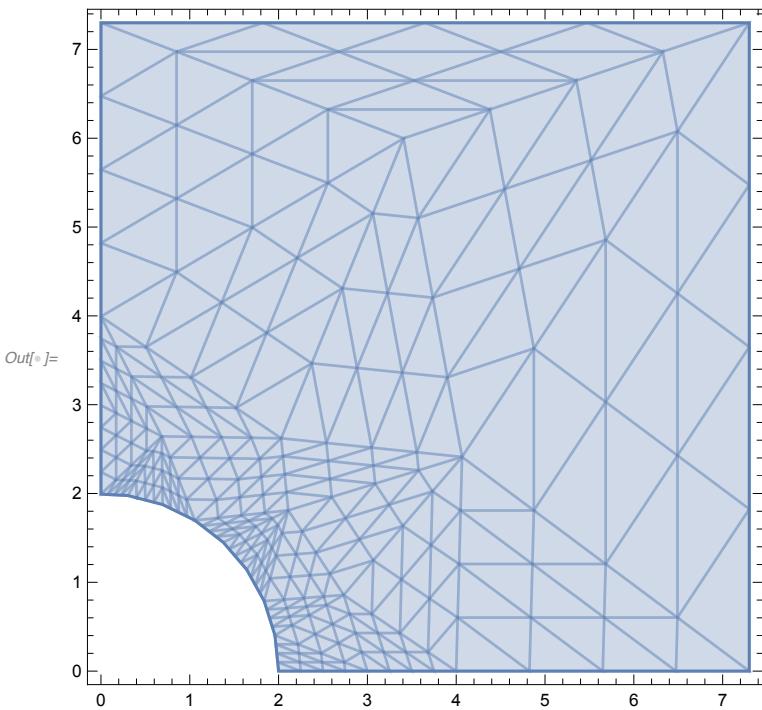
Out[=]/TraditionalForm=

$$\phi^{(0,2)}(r, z) + \frac{\phi^{(1,0)}(r, z)}{r} + \phi^{(2,0)}(r, z)$$


In[=]:= region = RegionDifference[Rectangle[{0, 0}, {b, h/2}], Disk[{0, 0}, a]]
Out[=]= BooleanRegion[!#1 && !#2 &, {Rectangle[{0, 0}, {7.3, 7.3}], Disk[{0, 0}, 2]}]

```

```
In[6]:= RegionPlot[region]
```



```
In[7]:= << NDSolve`FEM`
```

```
In[8]:= mesh = ToElementMesh[region, MaxCellMeasure -> 0.001]
```

```
Out[8]= ElementMesh[{{0., 7.3}, {0., 7.3}}, {TriangleElement[<79 083>]}]
```

```
In[9]:= Clear[z, r]
```

```
solc = NDSolveValue[
  {lpc == NeumannValue[0, r == 0 && a < z < b] + NeumannValue[0, z == 0 && a < r < b],
   DirichletCondition[\phi[r, z] == -1, r^2 + z^2 == a^2],
   \phi[r, h/2] == 0, \phi[b, z] == 0}, \phi, {r, z} \[Element] mesh
 ]
```

```
Out[9]= InterpolatingFunction[+  Domain: {{0., 7.3}, {0., 7.3}} ]
```

Data not in notebook. Store now

```
In[10]:= solc["ElementMesh"]
```

```
Out[10]= ElementMesh[{{0., 7.3}, {0., 7.3}}, {TriangleElement[<79 083>]}]
```

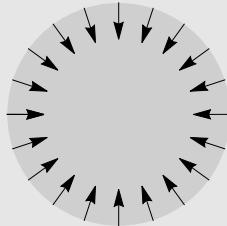
```

In[6]:= GetFromSolc[x_, z_] := If[Sqrt[x^2 + z^2] ≤ a, -1, solc[x, z]]
EstimateEField[x_, z_] := Module[{step, Er, Ez},
  step = 0.0005;
  Er =
    -(GetFromSolc[Abs[x + step / 2], Abs[z]] - GetFromSolc[Abs[x - step / 2], Abs[z]]) /
    step;
  Ez =
    -(GetFromSolc[Abs[x], Abs[z + step / 2]] - GetFromSolc[Abs[x], Abs[z - step / 2]]) /
    step;
  {Er, Ez}
]

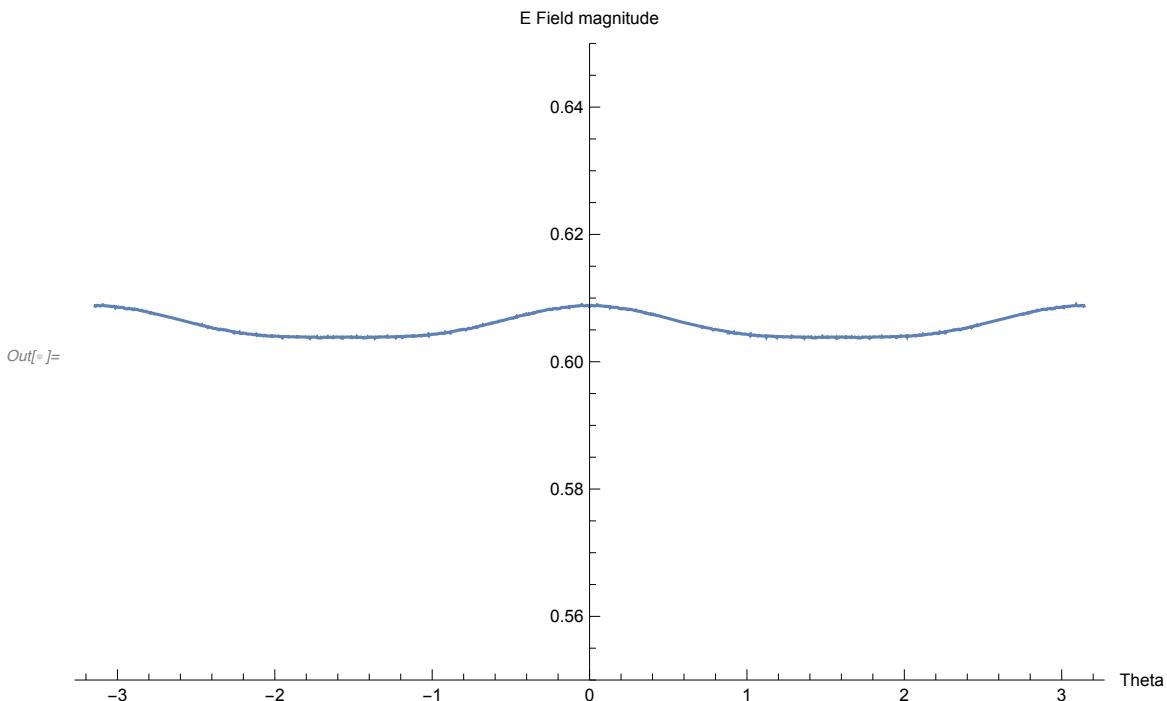
In[7]:= efieldvecs = Table[
  {{(a + gs) Cos[theta], (a + gs) Sin[theta]}, {(a + gs) Cos[theta], (a + gs) Sin[theta]} +
   {-GetEField[1, (a + gs) Cos[theta], 0, (a + gs) Sin[theta]][[1]],
    -GetEField[1, (a + gs) Cos[theta], 0, (a + gs) Sin[theta]][[3]]}},
  {theta, -Pi, Pi, Pi / 10}];
Show[Graphics[{Thickness[0.001], Arrowheads[0.02], Arrow[efieldvecs]}],
  Graphics[{Opacity[0.1], Disk[{0, 0}, a], Rectangle[{-b, -h / 2}, {b, h / 2}]}]]

```

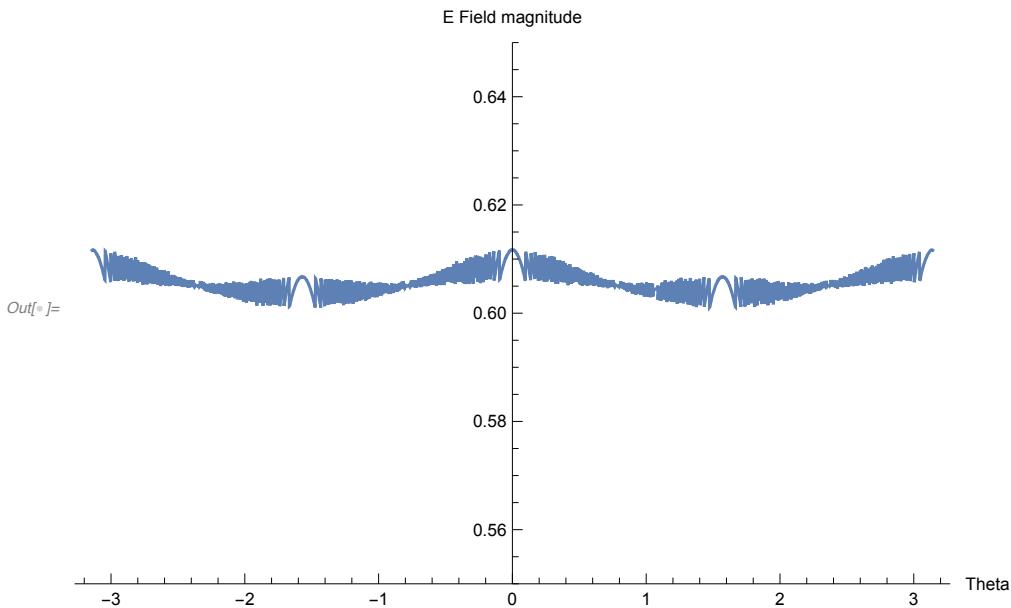
Out[7]=



```
In[6]:= d = 0.1;
Plot[Norm[EstimateEField[(a + d) Cos[theta], (a + d) Sin[theta]]], {theta, -Pi, Pi},
PlotRange -> {0.55, 0.65}, AxesLabel -> {"Theta", "E Field magnitude"}]
```



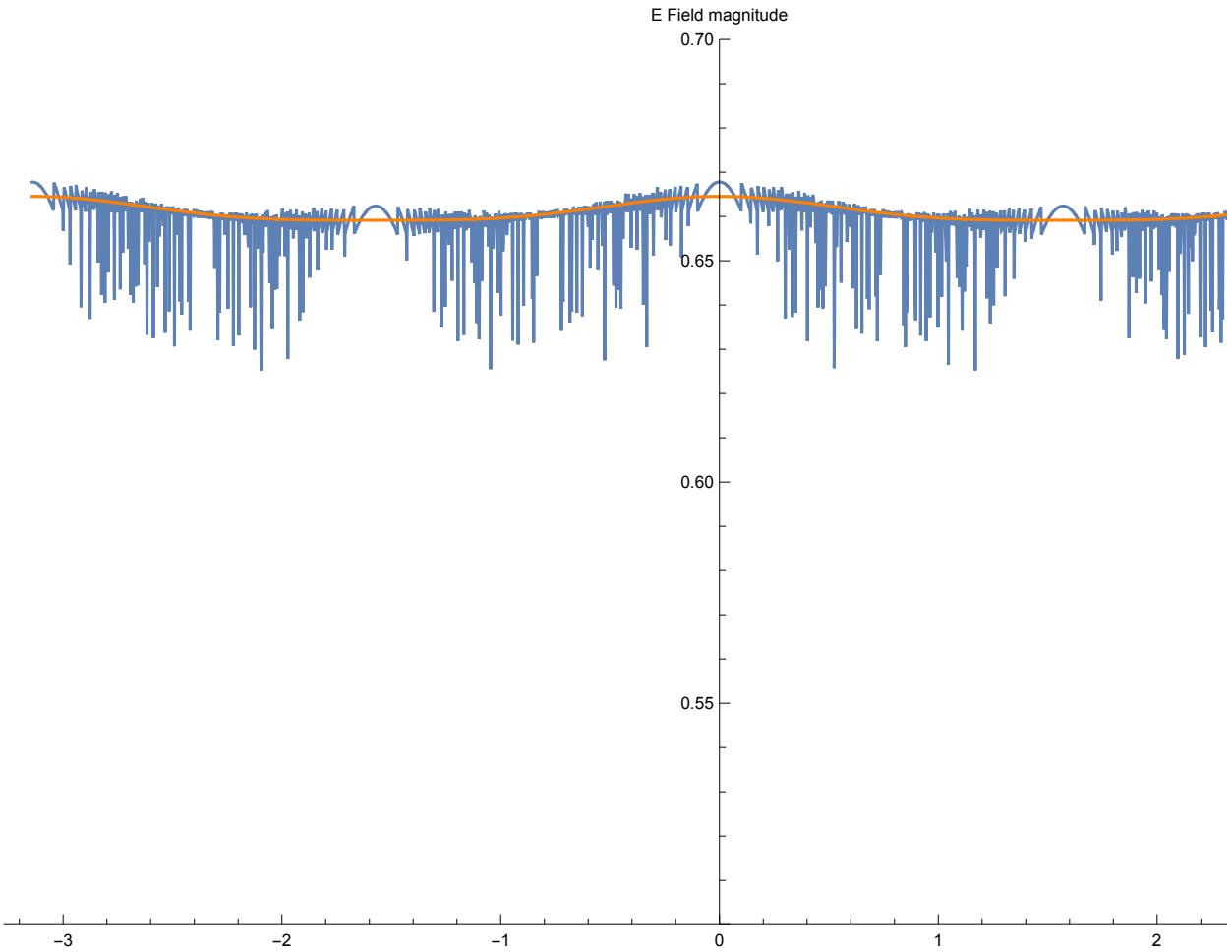
```
In[7]:= Plot[Norm[GetEField[1, (a + d) Cos[theta], 0, (a + d) Sin[theta]]], {theta, -Pi, Pi},
PlotRange -> {0.55, .65}, AxesLabel -> {"Theta", "E Field magnitude"}]
```



```
In[6]:= d = 1;
errors = Table[(Norm[EstimateEField[(a + d) Cos[theta], (a + d) Sin[theta]]] -
    Norm[GetEField[1, (a + d) Cos[theta], 0, (a + d) Sin[theta]]]) /
    Norm[EstimateEField[(a + d) Cos[theta], (a + d) Sin[theta]]],
{theta, -Pi, Pi, Pi / 20000}];
100 * Max[errors]

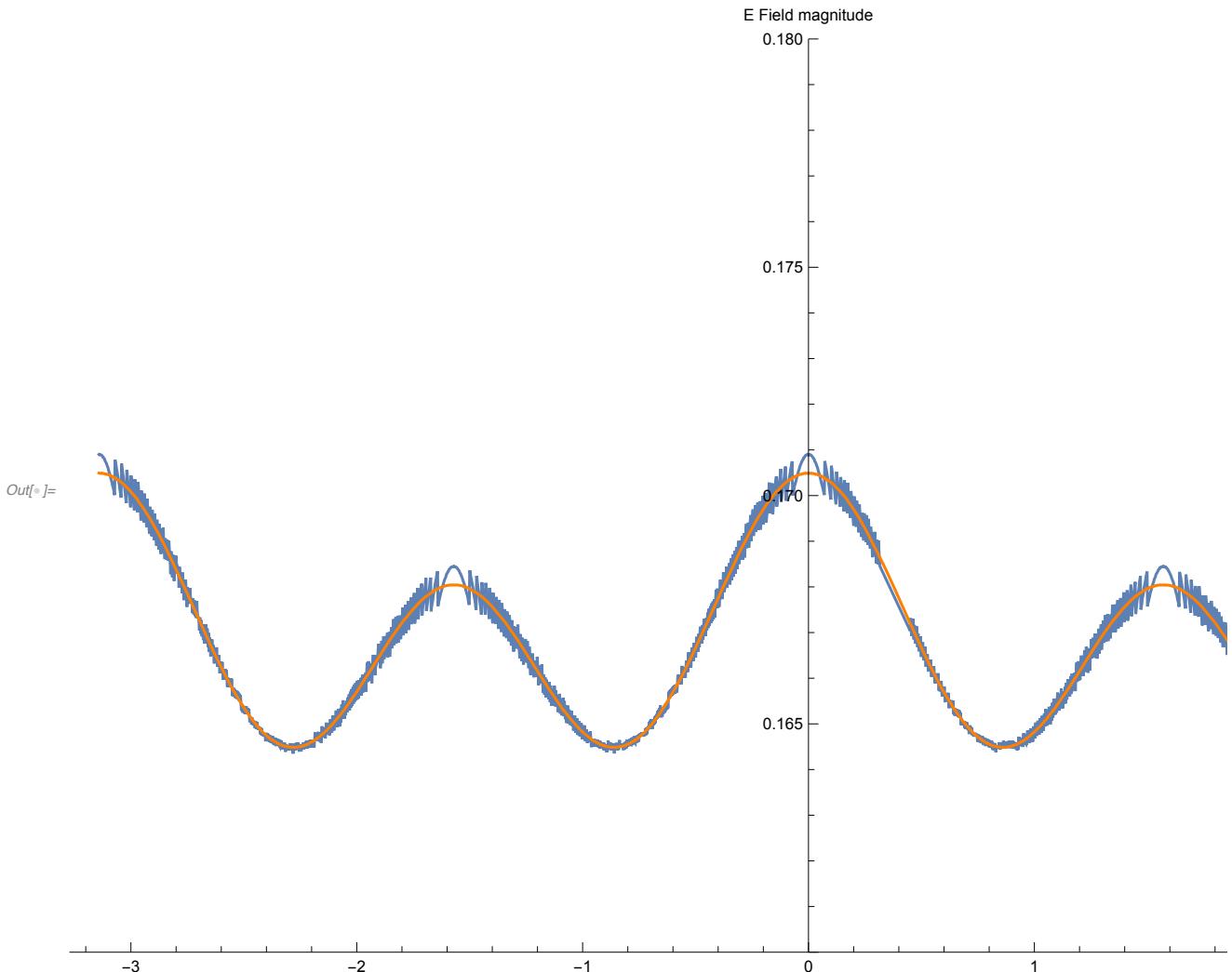
Out[6]= 0.32659

In[7]:= d = 0.01;
Show[Plot[Norm[GetEField[1, (a + d) Cos[theta], 0, (a + d) Sin[theta]]],
{theta, -Pi, Pi}, PlotRange -> {0.5, 0.7}],
Plot[Norm[EstimateEField[(a + d) Cos[theta], (a + d) Sin[theta]]], {theta, -Pi, Pi},
PlotStyle -> Orange], AxesLabel -> {"Theta", "E Field magnitude"}]
```

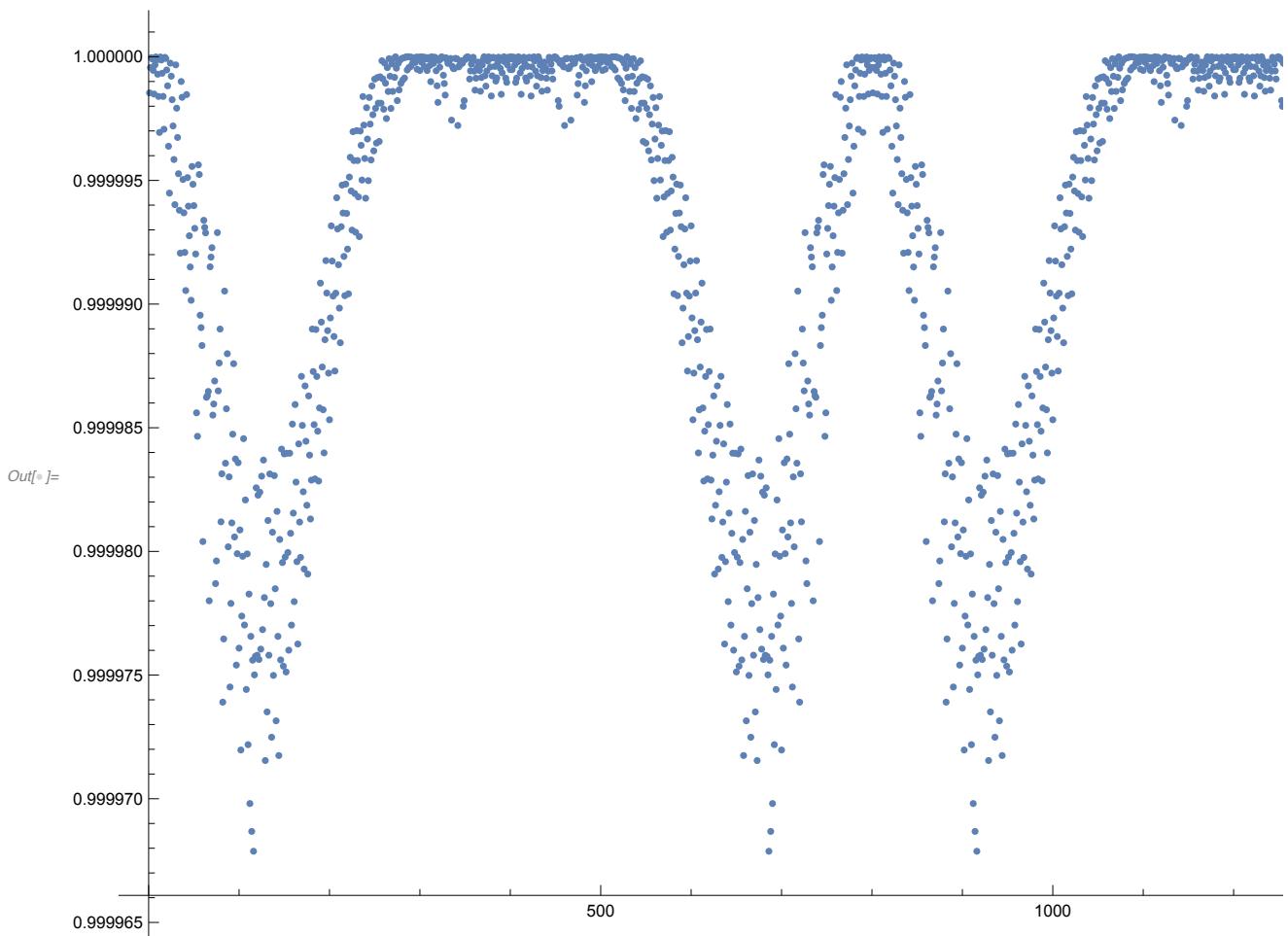


```
In[6]:= d = 2;
```

```
Show[Plot[Norm[GetEField[1, (a + d) Cos[theta], 0, (a + d) Sin[theta]]], {theta, -Pi, Pi}, PlotRange -> {0.16, 0.18}], Plot[Norm[EstimateEField[(a + d) Cos[theta], (a + d) Sin[theta]]], {theta, -Pi, Pi}, PlotStyle -> Orange], AxesLabel -> {"Theta", "E Field magnitude"}]
```

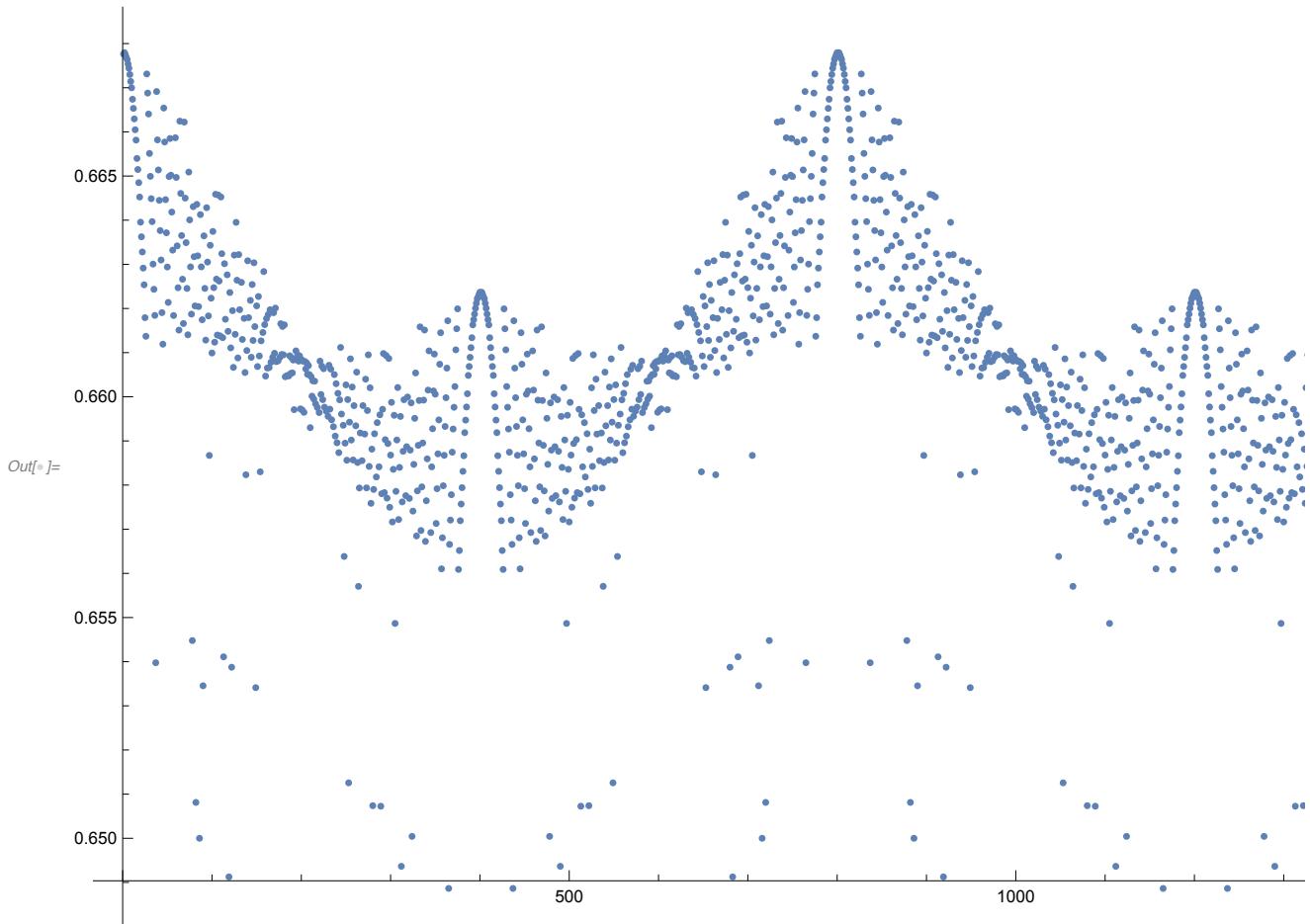


```
In[ $\circ$ ]:= d = 1
directions = Table[(GetEField[1, (a + d) Cos[theta], 0, (a + d) Sin[theta]].
{Cos[theta], 0, Sin[theta]})) /
Norm[GetEField[1, (a + d) Cos[theta], 0, (a + d) Sin[theta]]], {theta,
-Pi, Pi, 2 Pi / 1600}];
ListPlot[directions]
```

Out[\circ]= 1

```
In[ $\circ$ ]:= (*degrees of inaccuracy*)
ArcCos[Min[directions]] 180 / Pi
Out[ $\circ$ ]= 0.459241
```

```
In[6]:= mags = Table[Norm[GetEField[1, (a + gs) Cos[theta], 0, (a + gs) Sin[theta]]], {theta, -Pi, Pi, 2 Pi / 1600}];  
ListPlot[mags]
```



```
In[7]:= testPoints = {{4.608842669337616` , 6.563585696985424`},  
{3.707180200517526` , 2.176750103748627`},  
{3.792037566239333` , 4.787781408276944`}, {3.855778982477084` ,  
6.2237642461949205`}, {0.5816942256367047` , 6.431626072609152`},  
{1.866278784790517` , 1.5856972073106224`}, {2.905554076101517` ,  
0.3042719406065806`}, {4.575278216079155` , 4.368535882641796`},  
{0.15228864787249807` , 6.7965613936293`}, {3.0835338797819465` ,  
1.126015266510481`}, {0.8357538179378818` , 4.235467354793811`},  
{1.838948320701085` , 4.892962053921949`}, {0.6921786231648643` ,  
1.984597266088046`}, {1.6282948913230884` , 6.1271049030900615`},  
{6.260241423248773` , 4.049899219509415`}, {4.314099980656643` ,  
6.682551100251801`}, {3.13939323464767` , 3.2866671843639033`},  
{2.5535700266727304` , 6.458169956398989`}, {6.4600376108498825` ,  
2.845434570939234`}, {5.915655343265746` , 5.078591103153024`}};
```

```

In[6]:= testResults = Table[
  Module[{e1, e2, n1, n2, diff, cos},
    e1 = -{GetEField[1, point[[1]], 0, point[[2]]][1],
           GetEField[1, point[[1]], 0, point[[2]]][3]};
    e2 = EstimateEField[point[[1]], point[[2]]];
    n1 = Norm[e1];
    n2 = Norm[e2];
    diff = 2 (n1 - n2) / (n1 + n2);
    cos = e1.e2 / (n1 n2);
    {point, e1, e2, n1, n2, diff, cos}
  ], {point, testPoints}]];
testResults = Insert[testResults, {"Point", "E from File",
  "E from Laplace", "Norm 1", "Norm 2", "Difference", "Cos"}, 1];
Print[phiFile];
Grid[testResults, Frame → All, Spacings → 2]

```

Phi file: 2/5/2023 Phi from 2D cylindrical solve

Point	E from File	E from Laplace	Norm 1	Norm 2	Difference	Cos
{4.60884, 6.56359}	{-0.00958667, -0.0314003}	{-0.00958494, -0.0314021}	0.0328311	0.0328324	-0.0000379209	1.
{3.70718, 2.17675}	{-0.125929, -0.0692497}	{-0.125842, -0.0692609}	0.143714	0.143643	0.000492979	1.
{3.79204, 4.78778}	{-0.0390244, -0.0523992}	{-0.0390244, -0.052376}	0.0653344	0.0653158	0.00028492	1.
{3.85578, 6.22376}	{-0.0145572, -0.0426574}	{-0.0145568, -0.0426622}	0.0450728	0.0450773	-0.0000949782	1.
{0.581694, 6.43163}	{-0.0031235, -0.0775592}	{-0.00310674, -0.0775946}	0.0776221	0.0776567	-0.000446382	1.
{1.86628, 1.5857}	{-0.33956, -0.287458}	{-0.339769, -0.287402}	0.445199	0.44502	0.000403887	1.

	{2.90555, 0.304272 }	{-0.3138 46, -0.0323 101}	{-0.3137 43, -0.0322 276}	0.315505	0.315394	0.000350 693	1.
	{4.57528, 4.36854}	{-0.0442 936, -0.0390 171}	{-0.0442 909, -0.0390 052}	0.0590275	0.0590177	0.000166 391	1.
	{0.152289 , 6.79656}	{-0.0004 72751, -0.0755 912}	{-0.0004 64495, -0.0755 819}	0.0755926	0.0755833	0.000123 898	1.
Out[=]	{3.08353, 1.12602}	{-0.2346 41, -0.0833 52}	{-0.2348 26, -0.0833 882}	0.249006	0.249192	-0.00074 9257	1.
	{0.835754 , 4.23547}	{-0.0265 261, -0.1420 51}	{-0.0265 446, -0.1420 25}	0.144506	0.144484	0.000153 714	1.
	{1.83895, 4.89296}	{-0.0306 978, -0.0936 121}	{-0.0307 348, -0.0936 649}	0.0985169	0.0985786	-0.00062 5871	1.
	{0.692179 , 1.9846}	{-0.19912 , -0.5690 07}	{-0.1985 92, -0.5692 15}	0.602842	0.602863	-0.00003 62365	1.
	{1.62829, 6.1271}	{-0.0109 15, -0.0733 952}	{-0.0109 32, -0.0733 701}	0.0742024	0.07418	0.000301 489	1.
	{6.26024, 4.0499}	{-0.0388 513, -0.0123 728}	{-0.0388 805, -0.0123 718}	0.0407739	0.0408013	-0.00067 2549	1.
	{4.3141, 6.68255}	{-0.0081 1807, -0.0351 415}	{-0.0081 1806, -0.0351 454}	0.036067	0.0360708	-0.00010 719	1.

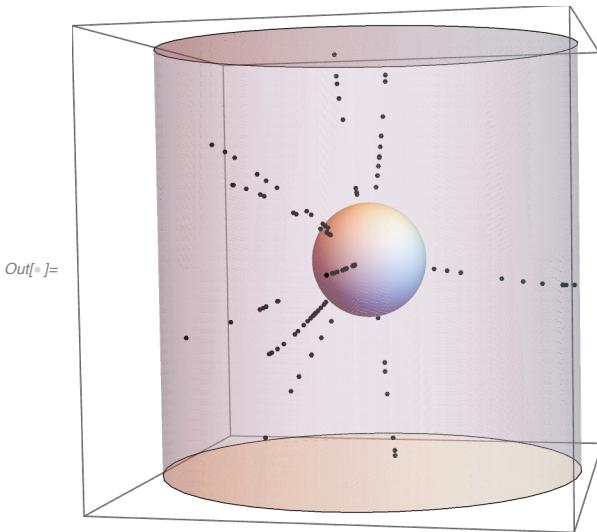
{3.13939, 3.28667}	{-0.0881, 766, -0.0900, 567}	{-0.0881, 195, -0.0900, 236}	0.126037	0.125973	0.000505 233	1.
{2.55357, 6.45817}	{-0.0101, 184, -0.0595, 603}	{-0.0101, 166, -0.0595, 414}	0.0604137	0.0603947	0.000314 094	1.
{6.46004, 2.84543}	{-0.0527, 552, -0.0092, 6054}	{-0.0528, 024, -0.0092, 6173}	0.0535619	0.0536085	-0.00087 0031	1.
{5.91566, 5.07859}	{-0.0269, 175, -0.0162, 529}	{-0.0269, 149, -0.0162, 49}	0.0314437	0.0314395	0.000132 705	1.

```
In[~]:= testResultsv2 = Table[{  
    point,  
    -GetPotential[1, point[[1]], 0, point[[2]]],  
    solc[point[[1]], point[[2]]],  
    100 * (solc[point[[1]], point[[2]]] + GetPotential[1, point[[1]], 0, point[[2]]]) /  
        solc[point[[1]], point[[2]]]} // N,  
    {point, testPoints}];  
testResultsv2 = Insert[testResultsv2, {"Point",  
    "Potential from File", "Potential from Laplace", "Percent Error"}, 1];  
Grid[testResultsv2, Frame → All, Spacings → 2]
```

Point	Potential from File	Potential from Laplace	Percent Error
{4.60884, 6.56359}	-0.0228437	-0.0228436	-0.0000893453
{3.70718, 2.17675}	-0.285638	-0.285638	-0.000119639
{3.79204, 4.78778}	-0.113569	-0.113568	-0.0000966593
{3.85578, 6.22376}	-0.0444733	-0.0444733	-0.0000920436
{0.581694, 6.43163}	-0.0647829	-0.0647828	-0.0000635923
{1.86628, 1.5857}	-0.755247	-0.755245	-0.000283062
{2.90555, 0.304272}	-0.576356	-0.576355	-0.000247503
{4.57528, 4.36854}	-0.0999098	-0.0999097	-0.000092579
{0.152289, 6.79656}	-0.0375394	-0.0375394	-0.000113133
{3.08353, 1.12602}	-0.476133	-0.476132	-0.000172861
{0.835754, 4.23547}	-0.28458	-0.284579	-0.000140282
{1.83895, 4.89296}	-0.177993	-0.177992	-0.000109267
{0.692179, 1.9846}	-0.935483	-0.935479	-0.000463012
{1.62829, 6.1271}	-0.0807665	-0.0807664	-0.000110662
{6.26024, 4.0499}	-0.0374481	-0.0374481	-5.93771 × 10 ⁻⁶
{4.3141, 6.68255}	-0.0214964	-0.0214964	-0.0000969176
{3.13939, 3.28667}	-0.256277	-0.256277	-0.0000666832
{2.55357, 6.45817}	-0.0487486	-0.0487486	-0.0000427234
{6.46004, 2.84543}	-0.0414332	-0.0414332	0.0000257004
{5.91566, 5.07859}	-0.0340119	-0.0340119	-0.000136803

```
In[=]:= points = Reap[AvgElectronsEulerSC[53 / 100, 1, 15, 2 / 5, 99 / 100, True, 10]] [[2]]
1000 * RMS travel error: 1.77026
V/NC 1.5
λ: 53
      100
Geometry: SC
Collisions: 10.0 ± 1.3
Mean path: 8.8186
```

```
In[6]:= Show[Graphics3D[{Opacity[0.3], Cylinder[{{0, 0, -7.3}, {0, 0, 7.3}}, 7.3]}, 
Graphics3D[Ball[{0, 0, 0}, 2]], 
Graphics3D[{Black, PointSize[0.01], Point[points[[1]]]}]]]
```



Sphere-in-cylinder MC's

```
In[7]:= AvgElectronsEulerSC[\lambda_, Ni_, Ui_, dt_, tbound_, print_, getbins_, count_] :=
Module[{eErrorList, V, result, stack, collisions, paths, svals, countbin,
electronsbin, squaredbin, lbin, lbinSquared, data, seedcount, totalpaths, Nc},
V = Ui * Ni;
totalpaths = 0;
seedcount = 1;
collisions = {};
eErrorList = {};
paths = {};
svals = 0;
If[getbins,
countbin = Table[0, 2 / binwidth];
electronsbin = Table[0, 2 / binwidth];
squaredbin = Table[0, 2 / binwidth];
lbin = Table[0, lmax / 2 + 1];
lbinSquared = Table[0, lmax / 2 + 1, lmax / 2 + 1]];
data = {};
stack = CreateDataStructure["Stack"];
result = Table[
Module[{x0, x1, y0, y1, z0, z1, numElectrons, cosθ, φ,
s, vx, vy, vz, energy, eError, numCollisions, pathval, sval, idx,
cθθ, φθ, SH, l, l2, collisionE, cost, phi, start, path, input},
numElectrons = 1;
```

```

numCollisions = 0;
stack["Push", GetStartingPos[]];
{cθ0, φ0} = GetAngles[x0, y0, z0];
While[! stack["EmptyQ"],
  {x0, y0, z0, vx, vy, vz} = stack["Pop"];
  cosθ = 1;
  φ = 0;
  energy = 0;
  path = 0;
  While[
    s = RandomVariate[ExponentialDistribution[1 / λ], WorkingPrecision → 16];
    {x1, y1, z1, vx, vy, vz, energy, eError, pathval, sval} =
      GetNewPositionsEuler[dt, s, V, x0, y0, z0, vx, vy, vz];
    AppendTo[eErrorList, eError];
    InBounds[x1, y1, z1],
    path += pathval;
    AppendTo[paths, pathval];
    svals += s;
    numCollisions++;
    If[energy ≥ Ui,
      numElectrons++;
      stack["Push", {x1, y1, z1, 0, 0, 0}];
      energy -= Ui,
      collisionE =
        SetPrecision[Ui * RandomVariate[UniformDistribution[{1 / 10, 1 / 2}], 16];
      energy -= collisionE;
      If[energy < 0, energy = 0];
    ];
    {vx, vy, vz, cost, phi} = UpdateVelocity[energy, vx, vy, vz, tbound];
    {x0, y0, z0} = {x1, y1, z1};
  ];
  totalpaths += path;
];
AppendTo[data, {cθ0, numElectrons}];
If[getbins,
  idx = FindBinNum[cθ0, binwidth];
  countbin[[idx]] += 1;
  electronsbin[[idx]] += numElectrons;
  squaredbin[[idx]] += numElectrons^2;
  SH = numElectrons Table[
    Conjugate[SphericalHarmonicY[l, 0, ArcCos[cθ0], φ0]], {l, 0, lmax, 2}];
  For[l = 0, l ≤ lmax, l += 2,
    lbin[[l / 2 + 1]] += SH[[l / 2 + 1]];
  For[l2 = 0, l2 ≤ lmax, l2 += 2,

```

```

lbinSquared[[l / 2 + 1, l2 / 2 + 1]] += SH[l / 2 + 1] × SH[l2 / 2 + 1]
]
];
];
AppendTo[collisions, numCollisions];
numElectrons,
{count}];
If[print,
Print["1000 * RMS travel error: ", 1000 * RootMeanSquare[eErrorList]];
Print["V/NC ", Ni Ui λ / 5.3];
Print["λ: ", λ];
Print["Geometry: ", ID];
Print["Collisions: ",
Around[Mean[collisions], StandardDeviation[collisions] / Sqrt[count]]];
Print["Mean path: ", totalpaths / count / λ];
Print[phiFile]
];
Sow[data ];
Around[Mean[result], StandardDeviation[result] / Sqrt[count]]]

```

Creating a faster version to run in-parallel

```

In[=]:= AbsoluteTiming[AvgElectronsEulerSCFAST[1 / 2, 10, 15, 2 / 5, 4 / 5, True, False, 100]]
1000 * RMS travel error: 14.6545
V/NC 14.1509
λ: 1
2
Geometry: SC
Collisions: 226. ± 10.
Mean path: 0
Phi file: 2/5/2023 Phi from 2D cylindrical solve
Out[=]= {115.108, 36.9 ± 1.4}

In[=]:= AvgElectronsEulerSCFAST[λ_, Ni_, Ui_, dt_, tbound_, print_, getbins_, count_] :=
Module[{masterdata, eErrorList, V, result, stack, collisions,
paths, svals, eErrorListMASTER, countbin, electronsbin,
squaredbin, lbin, lbinSquared, seedcount, totalpaths, Nc},
V = Ui * Ni;
seedcount = 1;
collisions = {};
(*each kernel creates its own version of eErrorList*)
eErrorListMASTER = {};
paths = {}];

```

```

svals = 0;
If[getbins,
  countbin = Table[0, 2 / binwidth];
  electronsbin = Table[0, 2 / binwidth];
  squaredbin = Table[0, 2 / binwidth];
  lbin = Table[0, lmax / 2 + 1];
  lbinSquared = Table[0, lmax / 2 + 1, lmax / 2 + 1]];
masterdata = {};
result = ParallelTable[
  Module[{x0, x1, y0, y1, z0, z1, numElectrons, cosθ, φ, s, vx,
    vy, vz, energy, data, eError, numCollisions, pathval, sval, idx,
    cθ0, φ0, SH, l, l2, collisionE, cost, phi, start, path, input},
    numElectrons = 1;
    numCollisions = 0;
    eErrorList = {};
    stack = CreateDataStructure["Stack"];
    stack["Push", GetStartingPos[]];
    {cθ0, φ0} = GetAngles[x0, y0, z0];
    While[! stack["EmptyQ"],
      {x0, y0, z0, vx, vy, vz} = stack["Pop"];
      cosθ = 1;
      φ = 0;
      energy = 0;
      data = {};
      path = 0;
      While[
        s = RandomVariate[ExponentialDistribution[1 / λ], WorkingPrecision → 16];
        {x1, y1, z1, vx, vy, vz, energy, eError, pathval, sval} =
          GetNewPositionsEuler[dt, s, V, x0, y0, z0, vx, vy, vz];
        AppendTo[eErrorList, eError];
        InBounds[x1, y1, z1],
        path += pathval;
        AppendTo[paths, pathval];
        svals += s;
        numCollisions++;
        If[energy ≥ Ui,
          numElectrons++;
          stack["Push", {x1, y1, z1, 0, 0, 0}];
          energy -= Ui,
          collisionE =
            SetPrecision[Ui * RandomVariate[UniformDistribution[{1 / 10, 1 / 2}], 16];
          energy -= collisionE;
          If[energy < 0, energy = 0];
        ];
      ];
    ];
  ];
]

```

```

{vx, vy, vz, cost, phi} = UpdateVelocity[energy, vx, vy, vz, tbound];
{x0, y0, z0} = {x1, y1, z1};
];
];
AppendTo[data, {cθ0, numElectrons}];
If[getbins,
  idx = FindBinNum[cθ0, binwidth];
  countbin[[idx]] += 1;
  electronsbin[[idx]] += numElectrons;
  squaredbin[[idx]] += numElectrons^2;
  SH = numElectrons Table[
    Conjugate[SphericalHarmonicY[l, 0, ArcCos[cθ0], φ0]], {l, 0, lmax, 2}];
  For[l = 0, l ≤ lmax, l += 2,
    lbin[[l / 2 + 1]] += SH[[l / 2 + 1]];
    For[l2 = 0, l2 ≤ lmax, l2 += 2,
      lbinSquared[[l / 2 + 1, l2 / 2 + 1]] += SH[[l / 2 + 1]] × SH[[l2 / 2 + 1]]
    ]
  ];
];
{numElectrons, eErrorList, numCollisions, Last[data]}],
{count}];
collisions = Flatten[#[[3]] & /@ result];
eErrorListMASTER = Flatten[#[[2]] & /@ result];
masterdata = #[[4]] & /@ result;
result = Flatten[#[[1]] & /@ result];
If[print,
  Print["1000 * RMS travel error: ",
    1000 * RootMeanSquare[Flatten[eErrorListMASTER]]];
  Print["V/NC ", Ni U i λ / 5.3];
  Print["λ: ", λ];
  Print["Geometry: ", ID];
  Print["Collisions: ",
    Around[Mean[collisions], StandardDeviation[collisions] / Sqrt[count]]];
  Print[phiFile]
];
Sow[masterdata];
Around[Mean[result], StandardDeviation[result] / Sqrt[count]]]

```

Playing with spherical harmonics (using standardized testing data)

In[=]:= Import["/Users/wbowers/Downloads/testData1.txt", "Table"]

```

{ {Command, line:, emc, Ni=11.3, Nc=10, Ui=15, reps=10000, dt=1e-3, scatter=0.8,
  seed=1, shape=scf, phifile=cwPhi.txt, showpath=true, pathnum=-2},
  {}, {-0.24765, 121}, {0.58485, 30}, {0.82089, 64}, {0.51459, 60},
  {0.53806, 47}, {-0.75322, 74}, {-0.39226, 42}, {-0.80142, 69}, {0.51232, 13},
  {-0.17224, 62}, {-0.15049, 37}, {-0.57335, 67}, {0.66387, 29},
  {0.95469, 35}, {0.29807, 100}, ... 9968 ..., {0.25002, 37}, {-0.98531, 44},
  {-0.48698, 34}, {0.99284, 27}, {0.27253, 84}, {-0.37638, 49}, {0.43035, 20},
  {-0.37275, 37}, {-0.36154, 33}, {-0.33106, 32}, {0.40546, 16}, {0.35234, 68},
  {-0.67814, 12}, {0.12916, 37}, {-0.18555, 14}, {-0.93947, 85}, {0.26674, 68} }

Out[=]

```

large output | show less | show more | show all | set si | ze limit...

```

In[=]:= SHtestdata = Drop[Import["/Users/wbowers/Downloads/testData1.txt", "Table"], 2];
SHtestdata2 = Drop[Import["/Users/wbowers/Downloads/testData2.txt", "Table"], 2];

In[=]:= MakeElectronsPlot[data_] :=
  Module[{tenthOrderFit, SHYs, max, error, res, electrons, electronsSquared,
    counts, finalbins, half, temp, symbins, symcounts,
    symsquaredbins, errors, finalsymbins, cθ0, numelectrons},
    tenthOrderFit = LinearModelFit[data,
      Table[SphericalHarmonicY[l, 0, t, 0], {l, 0, 10, 2}] /. Cos[t] → z,
      z, IncludeConstantBasis → False];
    SHYs = Table[SphericalHarmonicY[l, 0, t, 0], {l, 0, lmax, 2}] /. {Cos[t] → z};
    max = NMaximize[{tenthOrderFit[z], -1 ≤ z ≤ 1}, z];
    error =
      Sqrt[SHYs.tenthOrderFit["CovarianceMatrix"].Transpose[SHYs]] /. max[[2]];
    res = Around[max[[1]], error];
    electrons = Table[0, 2/binwidth];
    electronsSquared = Table[0, 2/binwidth];
    counts = Table[0, 2/binwidth];
    For[i = 1, i ≤ Length[data], i++,
      {cθ0, numelectrons} = data[[i]];
      idx = FindBinNum[cθ0, binwidth];
      electrons[[idx]] += numelectrons;
      electronsSquared[[idx]] += numelectrons^2;
      counts[[idx]] += 1
    ];
    finalbins = electrons / counts;
    finalbins =
      Transpose[{Table[theta, {theta, -0.975, 0.975, binwidth}], finalbins}];

    half = Length[finalbins] / 2;
    temp = #[[2]] & /@ finalbins;
  ]

```

```

symbins = (temp[[1 ;; half]] + Reverse[temp[[half + 1 ;; Length[temp]]]]) / 2;
symcounts = (counts[[1 ;; half]] + Reverse[counts[[half + 1 ;; Length[counts]]]]) / 2;
symsquaredbins = (electronsSquared[[1 ;; half]] +
    Reverse[electronsSquared[[half + 1 ;; Length[electronsSquared]]]]) / 2;

symbins = Flatten[AppendTo[symbins, Reverse[symbins]]];
symcounts = Flatten[AppendTo[symcounts, Reverse[symcounts]]];
symsquaredbins = Flatten[AppendTo[symsquaredbins, Reverse[symsquaredbins]]];

symbins = Transpose[{Table[theta, {theta, -0.975, 0.975, binwidth}], symbins}];
symsquaredbins =
Transpose[{Table[theta, {theta, -0.975, 0.975, binwidth}], symsquaredbins}];

errors = Sqrt[(#[[2]] & /@ symsquaredbins) / symcounts - (#[[2]] & /@ symbins)^2] /
Sqrt[symcounts];
finalsymbins = symbins;

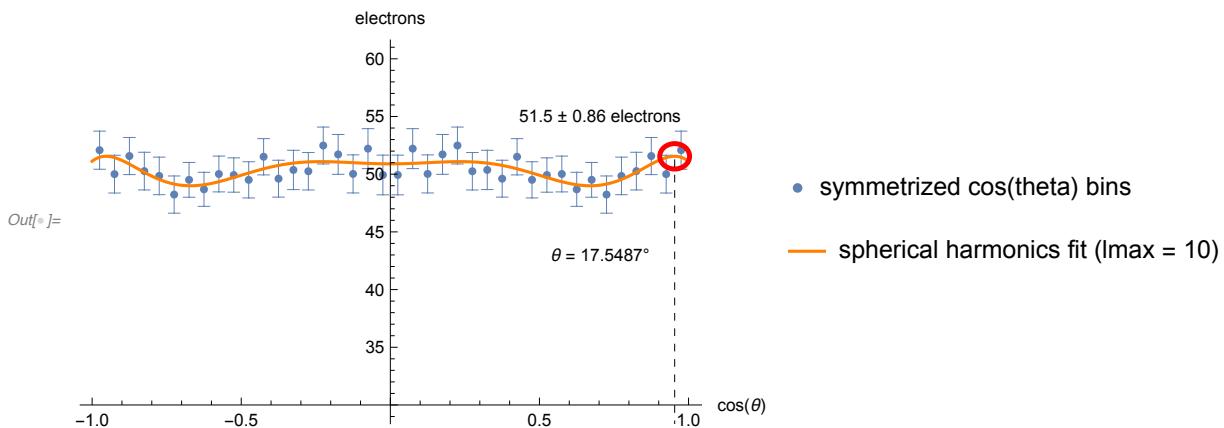
Print[finalbins];

For[i = 1, i ≤ Length[errors], i++,
finalsymbins[[i]] = {symbins[[i, 1]], Around[symbins[[i, 2]], errors[[i]]]}];
Show[ListPlot[finalsymbins, PlotLegends → {"symmetrized cos(theta) bins"}, ,
Plot[tenthOrderFit[x], {x, -1, 1}, PlotStyle → Orange,
PlotLegends → {"spherical harmonics fit (lmax = 10)"}, ,
Graphics[{Red, Thickness[0.008], Circle[{z /. max[[2]], max[[1]]}, {0.05, 1.1}]}], ,
Graphics[{Text[ToString[Round[res[[1]], 0.1]] <> " ± " <>
ToString[Round[res[[2]], 0.01]] <> " electrons", {(z /. max[[2]]) - 0.25, 55}], ,
Text["θ = " <> ToString[ArcCos[(z /. max[[2]]) * 1/Degree] <> "°",
{(z /. max[[2]]) - 0.25, 43}]}], ,
Graphics[{Dashed, Black, Line[{{z /. max[[2]], 20}, {z /. max[[2]], max[[1]]}}]}], ,
AxesOrigin → {0, 30}, PlotRange → {30, 60}, AxesLabel → {"cos(θ)", "electrons"}]
]
]

In[=]:= fig6 = MakeElectronsPlot[SHtestdata2]

```

$$\left\{ \left\{ -0.975, \frac{13213}{251} \right\}, \left\{ -0.925, \frac{12707}{249} \right\}, \left\{ -0.875, \frac{13771}{270} \right\}, \left\{ -0.825, \frac{13102}{257} \right\}, \left\{ -0.775, \frac{5949}{124} \right\}, \right. \\ \left. \left\{ -0.725, \frac{11422}{243} \right\}, \left\{ -0.675, \frac{11583}{226} \right\}, \left\{ -0.625, \frac{6272}{137} \right\}, \left\{ -0.575, \frac{13790}{269} \right\}, \left\{ -0.525, \frac{7031}{143} \right\}, \right. \\ \left. \left\{ -0.475, \frac{13244}{251} \right\}, \left\{ -0.425, \frac{667}{13} \right\}, \left\{ -0.375, \frac{12921}{253} \right\}, \left\{ -0.325, \frac{12401}{240} \right\}, \left\{ -0.275, \frac{1807}{35} \right\}, \right. \\ \left. \left\{ -0.225, \frac{12193}{244} \right\}, \left\{ -0.175, \frac{417}{8} \right\}, \left\{ -0.125, \frac{4049}{79} \right\}, \left\{ -0.075, \frac{12774}{241} \right\}, \left\{ -0.025, \frac{12330}{253} \right\}, \right. \\ \left. \left\{ 0.025, \frac{6851}{134} \right\}, \left\{ 0.075, \frac{11982}{233} \right\}, \left\{ 0.125, \frac{6197}{127} \right\}, \left\{ 0.175, \frac{12313}{240} \right\}, \left\{ 0.225, \frac{15179}{276} \right\}, \right. \\ \left. \left\{ 0.275, \frac{3275}{67} \right\}, \left\{ 0.325, \frac{11923}{243} \right\}, \left\{ 0.375, \frac{11893}{247} \right\}, \left\{ 0.425, \frac{6464}{125} \right\}, \left\{ 0.475, \frac{5273}{114} \right\}, \right. \\ \left. \left\{ 0.525, \frac{4207}{83} \right\}, \left\{ 0.575, \frac{11851}{243} \right\}, \left\{ 0.625, \frac{13570}{263} \right\}, \left\{ 0.675, \frac{12372}{259} \right\}, \left\{ 0.725, \frac{11673}{236} \right\}, \right. \\ \left. \left\{ 0.775, \frac{12621}{244} \right\}, \left\{ 0.825, \frac{11744}{237} \right\}, \left\{ 0.875, \frac{12879}{247} \right\}, \left\{ 0.925, \frac{10726}{219} \right\}, \left\{ 0.975, \frac{1391}{27} \right\} \right\}$$



1000 * RMS travel error: 149.96

V/NC 176.471

$$\lambda: \frac{53}{34}$$

Geometry: SC

Collisions: 42.47 ± 0.20

Phi file: 2/5/2023 Phi from 2D cylindrical solve

Out[6]=

```
{32.92 ± 0.15, {{{{0.360795, 71}, {0.598697, 88}, {-0.974832, 38}, {-0.767807, 18}, {-0.489686, 17}, {0.7662, 20}, {0.151377, 6}, {0.417244, 12}, {-0.205477, 45}, {-0.0468744, 14}, {-0.109721, 19}, {-0.530765, 50}, ... 49 977 ...}, {0.275548, 17}, {0.411143, 40}, {0.138401, 5}, {0.791972, 11}, {-0.960341, 7}, {0.774839, 4}, {0.908373, 22}, {0.861046, 62}, {-0.448823, 5}, {-0.865586, 27}, {-0.119944, 21}}}}}}
```

large output

show less

show more

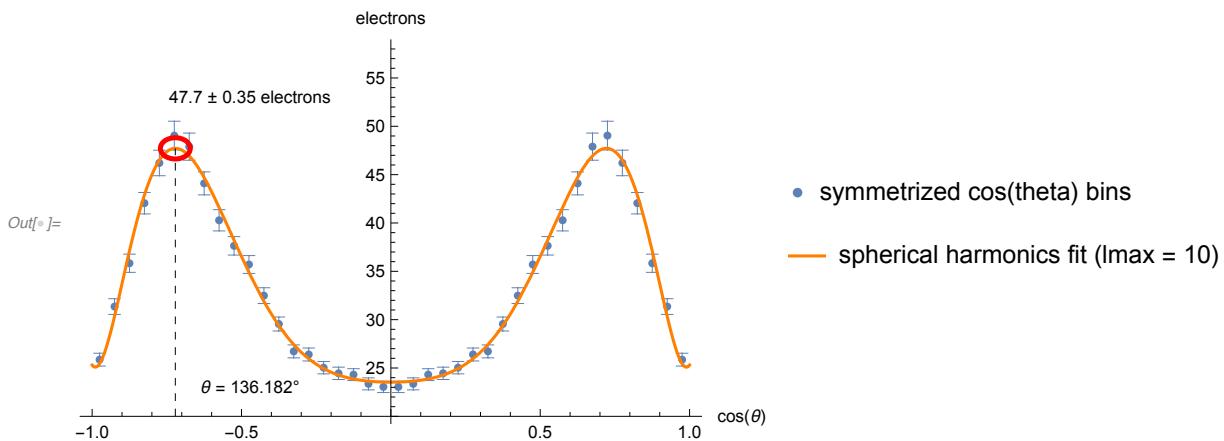
show all

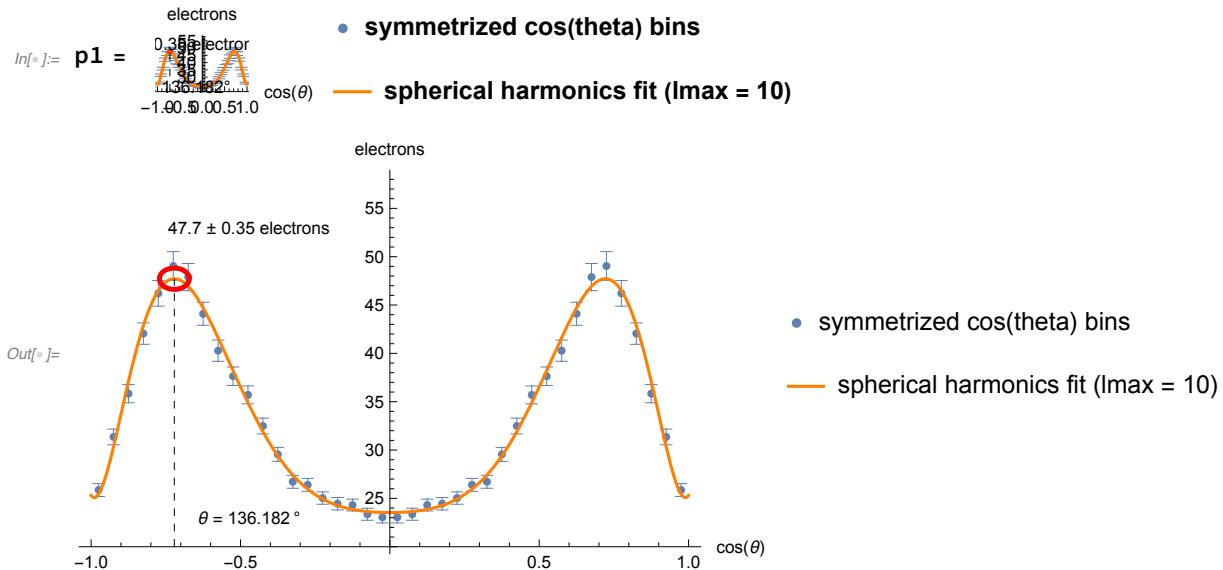
set si

ze limit...

In[6]:= MakeElectronsPlot[r[[2, 1, 1]]]

$$\left\{ \left\{ -0.975, \frac{103}{4} \right\}, \left\{ -0.925, \frac{42773}{1344} \right\}, \left\{ -0.875, \frac{15397}{421} \right\}, \left\{ -0.825, \frac{51929}{1243} \right\}, \left\{ -0.775, \frac{3556}{75} \right\}, \left\{ -0.725, \frac{49355}{1018} \right\}, \left\{ -0.675, \frac{8771}{178} \right\}, \left\{ -0.625, \frac{25888}{583} \right\}, \left\{ -0.575, \frac{49362}{1237} \right\}, \left\{ -0.525, \frac{25157}{652} \right\}, \left\{ -0.475, \frac{44791}{1250} \right\}, \left\{ -0.425, \frac{43137}{1366} \right\}, \left\{ -0.375, \frac{41359}{1373} \right\}, \left\{ -0.325, \frac{36101}{1330} \right\}, \left\{ -0.275, \frac{17747}{670} \right\}, \left\{ -0.225, \frac{10755}{443} \right\}, \left\{ -0.175, \frac{16970}{683} \right\}, \left\{ -0.125, \frac{4624}{193} \right\}, \left\{ -0.075, \frac{30052}{1285} \right\}, \left\{ -0.025, \frac{4984}{213} \right\}, \left\{ 0.025, \frac{29181}{1286} \right\}, \left\{ 0.075, \frac{28301}{1213} \right\}, \left\{ 0.125, \frac{16602}{673} \right\}, \left\{ 0.175, \frac{32179}{1338} \right\}, \left\{ 0.225, \frac{34483}{1337} \right\}, \left\{ 0.275, \frac{11339}{431} \right\}, \left\{ 0.325, \frac{34946}{1329} \right\}, \left\{ 0.375, \frac{39847}{1375} \right\}, \left\{ 0.425, \frac{10718}{321} \right\}, \left\{ 0.475, \frac{44224}{1243} \right\}, \left\{ 0.525, \frac{46561}{1269} \right\}, \left\{ 0.575, \frac{16627}{409} \right\}, \left\{ 0.625, \frac{17034}{389} \right\}, \left\{ 0.675, \frac{45491}{978} \right\}, \left\{ 0.725, \frac{3323}{67} \right\}, \left\{ 0.775, \frac{15707}{349} \right\}, \left\{ 0.825, \frac{24671}{583} \right\}, \left\{ 0.875, \frac{7651}{218} \right\}, \left\{ 0.925, \frac{20167}{653} \right\}, \left\{ 0.975, \frac{11404}{439} \right\} \}$$





In[2]:= $r2 = \text{Reap}[\text{AvgElectronsEulerSCFAST}[53 / 700, 171 / 10, 15, 2 / 5, 4 / 5, \text{True}, \text{False}, 5000]]$

1000 * RMS travel error: 21.1145

V/NC 3.66429

$$\lambda: \frac{53}{700}$$

Geometry: SC

Collisions: 2832. ± 14.

Phi file: 2/5/2023 Phi from 2D cylindrical solve

Out[2]:=

```
{45.12 ± 0.21, {{{{0.338099, 90}, {-0.962644, 19}, {-0.234118, 49}, {-0.608771, 79}, {-0.302092, 61}, {0.428077, 38}, {-0.63474, 21}, {-0.786352, 26}, {-0.811839, 34}, {0.378535, 49}, {0.752824, 27}, {-0.944244, 20}, ... 4977 ... , {-0.464546, 19}, {0.5626, 39}, {-0.0596975, 40}, {-0.438202, 33}, {-0.0369359, 67}, {0.045922, 79}, {0.339389, 71}, {-0.572022, 35}, {-0.210192, 58}, {-0.157178, 32}, {-0.0474263, 36}}}}}}
```

large output

show less

show more

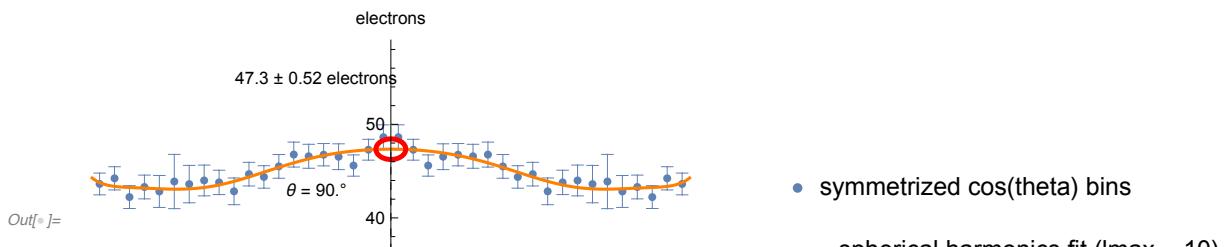
show all

set si

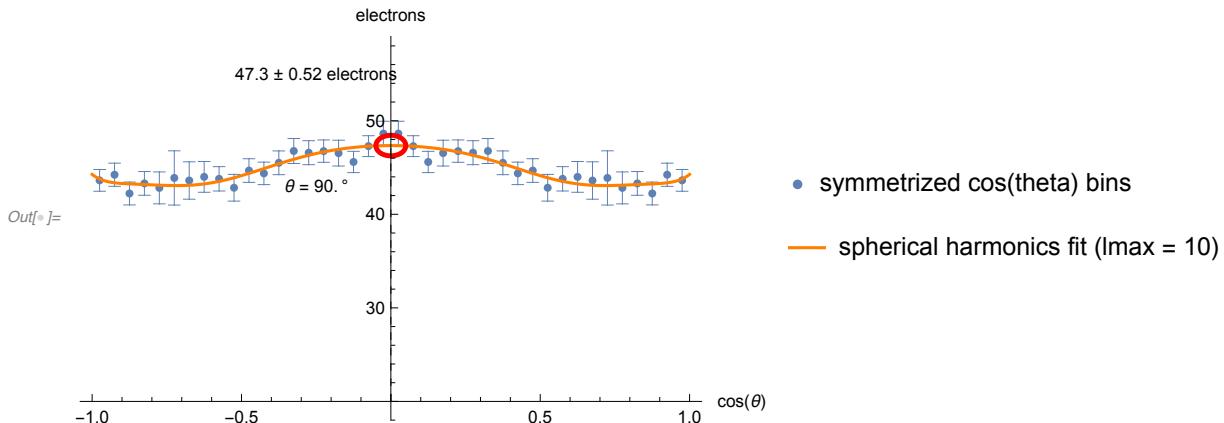
ze limit...

In[6]:= MakeElectronsPlot[r2[[2, 1, 1]]]

$$\left\{ \left\{ -0.975, \frac{897}{20} \right\}, \left\{ -0.925, \frac{2174}{49} \right\}, \left\{ -0.875, \frac{1171}{28} \right\}, \left\{ -0.825, \frac{4852}{113} \right\}, \left\{ -0.775, \frac{3752}{81} \right\}, \left\{ -0.725, \frac{1351}{30} \right\}, \left\{ -0.675, \frac{1840}{43} \right\}, \left\{ -0.625, \frac{2844}{65} \right\}, \left\{ -0.575, \frac{2237}{51} \right\}, \left\{ -0.525, \frac{1519}{36} \right\}, \left\{ -0.475, \frac{5539}{124} \right\}, \left\{ -0.425, \frac{6149}{137} \right\}, \left\{ -0.375, \frac{7432}{163} \right\}, \left\{ -0.325, \frac{6892}{149} \right\}, \left\{ -0.275, \frac{7209}{155} \right\}, \left\{ -0.225, \frac{6992}{147} \right\}, \left\{ -0.175, \frac{6577}{143} \right\}, \left\{ -0.125, \frac{7638}{163} \right\}, \left\{ -0.075, \frac{8171}{176} \right\}, \left\{ -0.025, \frac{3597}{73} \right\}, \left\{ 0.025, \frac{6957}{145} \right\}, \left\{ 0.075, \frac{7655}{159} \right\}, \left\{ 0.125, \frac{7403}{167} \right\}, \left\{ 0.175, \frac{7251}{154} \right\}, \left\{ 0.225, \frac{781}{17} \right\}, \left\{ 0.275, \frac{3033}{65} \right\}, \left\{ 0.325, \frac{6997}{148} \right\}, \left\{ 0.375, \frac{3179}{70} \right\}, \left\{ 0.425, \frac{6229}{142} \right\}, \left\{ 0.475, \frac{1832}{41} \right\}, \left\{ 0.525, \frac{1174}{27} \right\}, \left\{ 0.575, \frac{5378}{123} \right\}, \left\{ 0.625, \frac{3805}{86} \right\}, \left\{ 0.675, \frac{1644}{37} \right\}, \left\{ 0.725, \frac{983}{23} \right\}, \left\{ 0.775, \frac{3028}{77} \right\}, \left\{ 0.825, \frac{131}{3} \right\}, \left\{ 0.875, \frac{6607}{155} \right\}, \left\{ 0.925, \frac{6525}{148} \right\}, \left\{ 0.975, \frac{6318}{149} \right\} \right\}$$



In[7]:= p2 =



```
In[6]:= Export["/Users/wbowers/Documents/stem-fellowship/tif-figs/e-bins-1.TIFF",
  p1, ImageResolution → 600]
Export["/Users/wbowers/Documents/stem-fellowship/tif-figs/e-bins-2.TIFF",
  p2, ImageResolution → 600]

Out[6]= /Users/wbowers/Documents/stem-fellowship/tif-figs/e-bins-1.TIFF
Out[7]= /Users/wbowers/Documents/stem-fellowship/tif-figs/e-bins-2.TIFF
```

Using Spherical Harmonics to create max electrons function for sphere-in-cylinder

```
MaxElectrons[λ_, Ni_, reps_] :=
Module[{idx, list, i, thetaslist, constants, SHYs, fit, max, error},
  list = Last[
    Last[Reap[AvgElectronsEulerSC[λ, Ni, 15, 2/5, 4/5, False, True, reps]]][[1]]];
  fit = LinearModelFit[list,
    Table[SphericalHarmonicY[l, 0, theta, 0], {l, 0, lmax, 2}] /. Cos[theta] → z,
    z, IncludeConstantBasis → False];
  max = NMaximize[{fit[z], -1 ≤ z ≤ 1}, z];
  Print[z /. max[[2]]];
  SHYs =
    Table[SphericalHarmonicY[l, 0, theta, 0], {l, 0, lmax, 2}] /. Cos[theta] → z;
  error = Sqrt[SHYs.fit["CovarianceMatrix"].Transpose[SHYs]] /. max[[2]];
  Around[max[[1]], error]
]

In[7]:= MaxElectronsFAST[λ_, Ni_, reps_] :=
Module[{idx, list, i, thetaslist, constants, SHYs, fit, max, error},
  list = Last[Last[
    Reap[AvgElectronsEulerSCFAST[λ, Ni, 15, 2/5, 4/5, False, True, reps]]][[1]];
  Clear[theta, z];
  fit = LinearModelFit[list,
    Table[SphericalHarmonicY[l, 0, theta, 0], {l, 0, lmax, 2}] /. Cos[theta] → z,
    z, IncludeConstantBasis → False];
  max = NMaximize[{fit[z], -1 ≤ z ≤ 1}, z];
  Print[z /. max[[2]]];
  SHYs =
    Table[SphericalHarmonicY[l, 0, theta, 0], {l, 0, lmax, 2}] /. Cos[theta] → z;
  error = Sqrt[SHYs.fit["CovarianceMatrix"].Transpose[SHYs]] /. max[[2]];
  Around[max[[1]], error]
]
```

In[1]:= MaxElectronsFAST[53 / 100, 12.1, 500]

$$1.97348 \times 10^{-32}$$

Out[1]= 70. ± 4.

In[2]:= MaxElectronsFAST[53 / 100, 12.1, 500]

$$1.03398 \times 10^{-25}$$

Out[2]= 69. ± 4.

Ensuring exponential growth of ions for the parallel plate case (1D AND 3D)

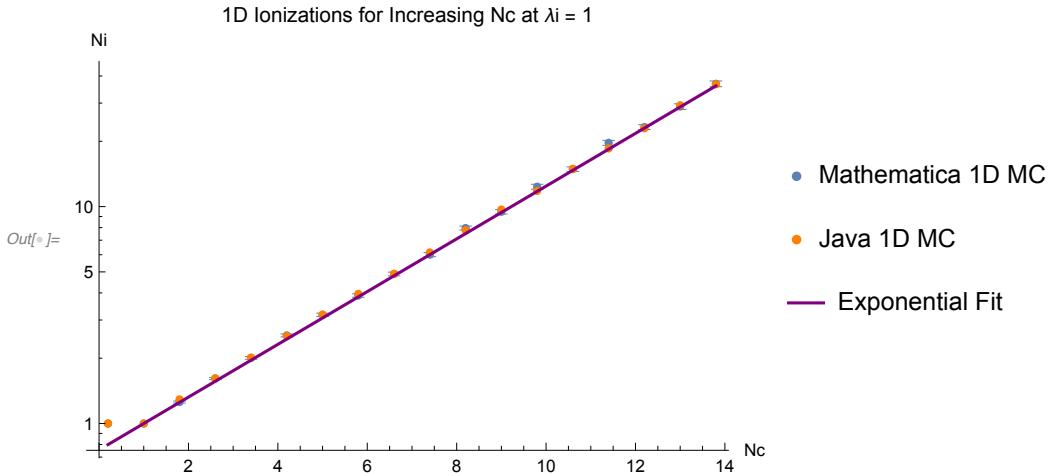
In[3]:= t1 = Table[{Nc, AvgElectronsWeighted[Nc, Nc, 1000]}, {Nc, 0.2, 14, 0.8}]

Out[3]= {{0.2, 1}, {1., 1}, {1.8, 1.256 ± 0.014}, {2.6, 1.612 ± 0.018}, {3.4, 2.006 ± 0.029}, {4.2, 2.55 ± 0.04}, {5., 3.17 ± 0.06}, {5.8, 3.89 ± 0.07}, {6.6, 4.89 ± 0.10}, {7.4, 6.00 ± 0.13}, {8.2, 7.96 ± 0.17}, {9., 9.47 ± 0.22}, {9.8, 12.35 ± 0.32}, {10.6, 14.9 ± 0.4}, {11.4, 19.7 ± 0.5}, {12.2, 23.3 ± 0.6}, {13., 28.9 ± 0.9}, {13.8, 36.8 ± 1.1}}

In[4]:= java1d = {{0.2, Round[1.0, 0.0]}, {1.0, Round[1.0, 0.0]}, {1.8, Round[1.292, 0.0]}, {2.6, Round[1.6152, 0.0]}, {3.4, Round[2.0128, 0.0]}, {4.2, Round[2.5308, 0.0]}, {5.0, Round[3.1744, 0.0]}, {5.8, Round[3.96, 0.0]}, {6.6, Round[4.8992, 0.0]}, {7.4, Round[6.146, 0.0]}, {8.2, Round[7.778, 0.0]}, {9.0, Round[9.6872, 0.0]}, {9.8, Round[11.818, 0.0]}, {10.6, Round[14.9372, 0.0]}, {11.4, Round[18.5404, 0.0]}, {12.2, Round[22.9588, 0.0]}, {13.0, Round[29.3304, 0.0]}, {13.8, Round[36.6168, 0.0]}, {14.6, Round[45.6408, 0.0]}}

Out[4]= {{0.2, 1.}, {1., 1.}, {1.8, 1.292}, {2.6, 1.6152}, {3.4, 2.0128}, {4.2, 2.5308}, {5., 3.1744}, {5.8, 3.96}, {6.6, 4.8992}, {7.4, 6.146}, {8.2, 7.778}, {9., 9.6872}, {9.8, 11.818}, {10.6, 14.9372}, {11.4, 18.5404}, {12.2, 22.9588}, {13., 29.3304}, {13.8, 36.6168}, {14.6, 45.6408}}

```
In[6]:= Show[ListLogPlot[t1, PlotLegends -> {"Mathematica 1D MC"}],  
ListLogPlot[java1d, PlotStyle -> Orange, PlotLegends -> {"Java 1D MC"}],  
LogPlot[Exp[0.28 (x - 1)], {x, 0.2, 13.8},  
PlotStyle -> Purple, PlotLegends -> {"Exponential Fit"}],  
PlotLabel -> "1D Ionizations for Increasing Nc at λi = 1", AxesLabel -> {Nc, Ni}]
```



```
Out[6]= t2 = Table[{Nc, AvgElectrons3D[Nc, Nc, 5, 15, 0.8, 100]}, {Nc, 0.2, 14, 0.8}]  
  
{ {0.2, 1}, {1., 1}, {1.8, 1.36 ± 0.05}, {2.6, 1.88 ± 0.07}, {3.4, 2.54 ± 0.10},  
{4.2, 3.41 ± 0.14}, {5., 5.31 ± 0.20}, {5.8, 6.94 ± 0.29}, {6.6, 10.15 ± 0.35},  
{7.4, 13.9 ± 0.5}, {8.2, 20.6 ± 0.7}, {9., 26.2 ± 1.0}, {9.8, 36.7 ± 1.1}, {10.6, 51.3 ± 1.7},  
{11.4, 74.4 ± 2.6}, {12.2, 102.1 ± 3.3}, {13., 136. ± 5.}, {13.8, 195. ± 7.} }
```

```
In[6]:= java3d = {{0.2, Around[1.0, 0.0]}, {1.0, Around[1.0, 0.0]}, {1.8, Around[1.3708, 0.009660380116744999]}, {2.6, Around[1.8896, 0.014252183552003757]}, {3.400000000000004, Around[2.6092, 0.01951076994892846]}, {4.2, Around[3.622, 0.02711690247797481]}, {5.0, Around[5.0172, 0.03788458346082171]}, {5.8, Around[7.038, 0.05230891319842165]}, {6.6, Around[9.6412, 0.071414459488258]}, {7.399999999999995, Around[13.5272, 0.0970370241918007]}, {8.2, Around[18.8308, 0.13820415530656086]}, {9.0, Around[26.5088, 0.18488258172148112]}, {9.8, Around[36.892, 0.25806691845333396]}, {10.600000000000001, Around[51.484, 0.353070726059242]}, {11.400000000000002, Around[71.3404, 0.49144933730344953]}, {12.200000000000003, Around[98.8472, 0.6953129517447517]}, {13.000000000000004, Around[139.1028, 0.961511358676537]}, {13.800000000000004, Around[194.0804, 1.3399496611201482]}}

Out[6]= {{0.2, 1.}, {1., 1.}, {1.8, 1.371 ± 0.010}, {2.6, 1.890 ± 0.014}, {3.4, 2.609 ± 0.020}, {4.2, 3.622 ± 0.027}, {5., 5.02 ± 0.04}, {5.8, 7.04 ± 0.05}, {6.6, 9.64 ± 0.07}, {7.4, 13.53 ± 0.10}, {8.2, 18.83 ± 0.14}, {9., 26.51 ± 0.18}, {9.8, 36.89 ± 0.26}, {10.6, 51.48 ± 0.35}, {11.4, 71.3 ± 0.5}, {12.2, 98.8 ± 0.7}, {13., 139.1 ± 1.0}, {13.8, 194.1 ± 1.3} }

In[7]:= Show[ListLogPlot[t2, PlotLegends → {"Mathematica 3D MC"}], ListLogPlot[java3d, PlotStyle → Orange, PlotLegends → {"Java 3D MC"}], LogPlot[Exp[0.41 (x - 1)], {x, 0.2, 13.8}, PlotStyle → Purple, PlotLegends → {"Exponential Fit"}], PlotLabel → "3D Ionizations for Increasing Nc at λi = 1", AxesLabel → {Nc, Ni}]

3D Ionizations for Increasing Nc at λi = 1
 $i_N$ 
 $c_N$ 

Out[7]=


```

Exploring 1D alphas

```
In[=]:= FindAlpha[λi_, reps_] := Module[{electrons, result,
  length, i, total, slopes, alpha, e, fittedmodel, a, b, c, d, params},
  electrons = Table[
    {Nc, Log[AvgElectronsWeighted[Nc, ((1/λi) Nc), reps]]}, {Nc, λi + 5, 40, 4}];
  fittedmodel = NonlinearModelFit[
    {#[[1]], #[[2, 1]]} & /@ electrons, a*x + b + c Exp[-d x], {a, b, c, d}, x];
  params = fittedmodel["ParameterTable"][[1, 1]];
  {Around[params[[2, 2]], params[[2, 3]]], Around[params[[3, 2]], params[[3, 3]]]}]

In[=]:= markovmodel[x_] := 1/x * ProductLog[x Exp[-x]]

In[=]:= javaalphas = {{0.25, Around[0.6602595438326849, 0.001904193660625115]}, {0.5, Around[0.47844948673235005, 0.001748144153630099]}, {0.75, Around[0.35821274297950323, 0.0031233087246661244]}, {1.0, Around[0.27433365451618424, 0.008691189348246631]}, {1.25, Around[0.21314895885656557, 0.013389317555812227]}, {1.5, Around[0.1669038047699682, 0.01749867647681158]}, {1.75, Around[0.13018046550569878, 0.014267346022973922]}, {2.0, Around[0.1015082557377325, 0.012894206875325626]}, {2.2500000000000004, Around[0.078742597634773, 0.011952013460590393]}, {2.5, Around[0.061520625296323746, 0.010962062286845231]}, {2.75, Around[0.04725434577446056, 0.00734274289001424]}, {3.0, Around[0.0358071249036958, 0.004959892432166247]}, {3.25, Around[0.02778741774482209, 0.004504199943380026]}, {3.5, Around[0.021572007247372436, 0.003796229153906858]}, {3.7500000000000004, Around[0.015726014436025514, 0.0022654561964774527]}, {4.0, Around[0.012178682642516597, 0.0017503399885901638]}, {4.25, Around[0.008830658060106817, 0.0010123646555449694]}}

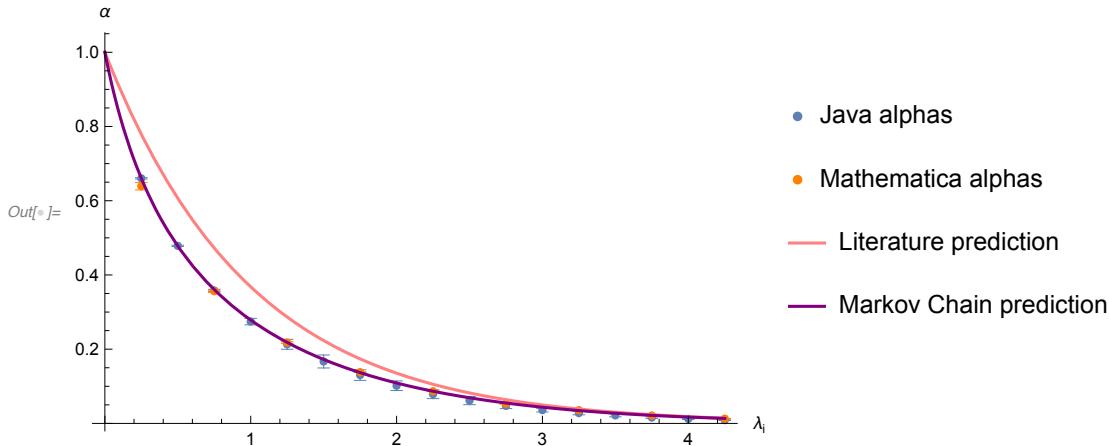
Out[=]= {{0.25, 0.6603 ± 0.0019}, {0.5, 0.4784 ± 0.0017}, {0.75, 0.3582 ± 0.0031}, {1., 0.274 ± 0.009}, {1.25, 0.213 ± 0.013}, {1.5, 0.167 ± 0.017}, {1.75, 0.130 ± 0.014}, {2., 0.102 ± 0.013}, {2.25, 0.079 ± 0.012}, {2.5, 0.062 ± 0.011}, {2.75, 0.047 ± 0.007}, {3., 0.036 ± 0.005}, {3.25, 0.028 ± 0.005}, {3.5, 0.022 ± 0.004}, {3.75, 0.0157 ± 0.0023}, {4., 0.0122 ± 0.0018}, {4.25, 0.0088 ± 0.0010} }

In[=]:= alphas1d = Table[{r, FindAlpha[r, 1000][[1]]}, {r, 0.25, 4.25, 0.5}]

Out[=]= {{0.25, 0.639 ± 0.010}, {0.75, 0.3563 ± 0.0028}, {1.25, 0.2179 ± 0.0022}, {1.75, 0.1382 ± 0.0009}, {2.25, 0.0868 ± 0.0009}, {2.75, 0.0529 ± 0.0011}, {3.25, 0.0356 ± 0.0018}, {3.75, 0.0212 ± 0.0006}, {4.25, 0.0129 ± 0.0006}}
```

```
In[6]:= fig8 = Show[ListPlot[javaalphas, PlotLegends -> {"Java alphas"}],
  ListPlot[alphas1d, PlotStyle -> Orange, PlotLegends -> {"Mathematica alphas"}],
  Plot[Exp[-x], {x, 0, 4.25}, PlotLegends -> {"Literature prediction"},
  PlotStyle -> Pink], Plot[1/x * ProductLog[x Exp[-x]], {x, 0, 4.25},
  PlotStyle -> Purple, PlotLegends -> {"Markov Chain prediction"}],
  PlotRange -> {0, 1}, AxesLabel -> {Subscript["λ", "i"], "α"},
  PlotLabel -> "1D Alphas with Markov Chain Model"]
```

1D Alphas with Markov Chain Model

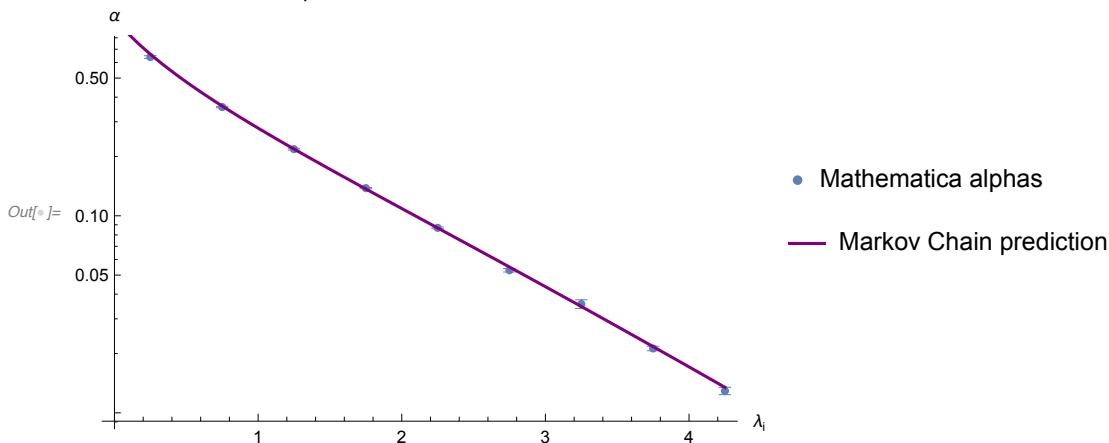


```
In[7]:= Export["/Users/wbowers/Documents/stem-fellowship/figures/alphas-plot.TIFF",
  fig8, ImageResolution -> 600]
```

```
Out[7]= /Users/wbowers/Documents/stem-fellowship/figures/alphas-plot.TIFF
```

```
In[8]:= Show[ListLogPlot[alphas1d, PlotLegends -> {"Mathematica alphas"}],
  LogPlot[1/x * ProductLog[x Exp[-x]], {x, 0, 4.25}, PlotStyle -> Purple,
  PlotLegends -> {"Markov Chain prediction"}], AxesLabel -> {Subscript["λ", "i"], "α"},
  PlotLabel -> "1D Alphas with Markov Chain Model"]
```

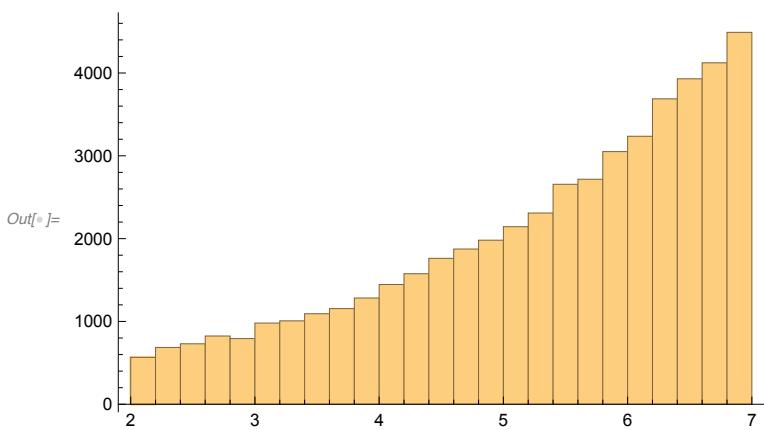
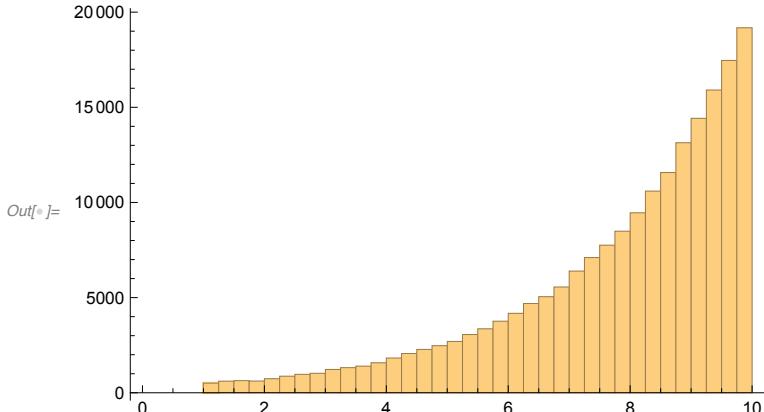
1D Alphas with Markov Chain Model



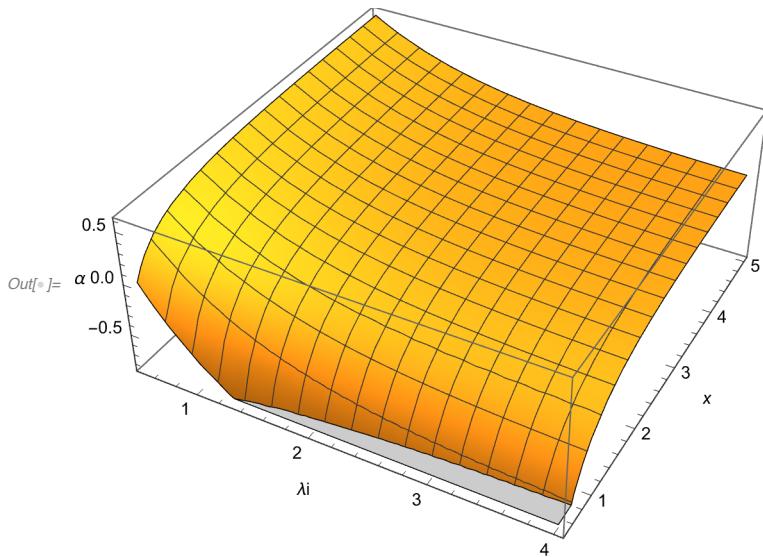
```
In[5]:= l1 = Reap[AvgElectrons3D[10, 10, 5000]]; l1[[1]]
Out[5]= 39.79 ± 0.19
```

Examining the sample space for the alpha calculation

```
In[6]:= Histogram[l1[[2, 1]], {0, 10, 0.25}]
Histogram[l1[[2, 1]], {2, 7, 0.2}]
```



```
In[6]:= Clear[\alpha]
FindAlphav2[\lambda i_, x_] := (Exp[-\lambda i] - Exp[-x]) / (1 - Exp[-x])
Plot3D[FindAlphav2[\lambda i, x], {\lambda i, 1/2, 4}, {x, 1/2, 5}, AxesLabel \rightarrow {\lambda i, x, \alpha}]
```



```
FindRoot[FindAlphav2[1/2, x] == FindAlpha[1/2][1], {x, 1, 6}]
FindRoot[FindAlphav2[1, x] == FindAlpha[1][1], {x, 1, 6}]
FindRoot[FindAlphav2[4, x] == FindAlpha[4][1], {x, 1, 6}]
```

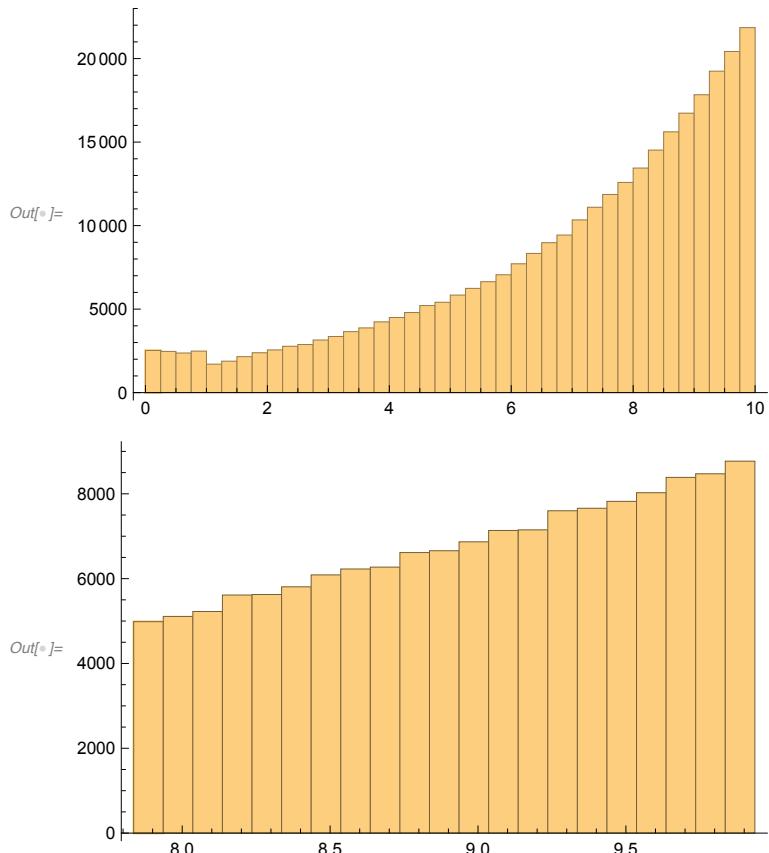
```
Out[6]= {x \rightarrow 1.42628}
```

```
Out[7]= {x \rightarrow 2.16402}
```

```
Out[8]= {x \rightarrow 6.78437}
```

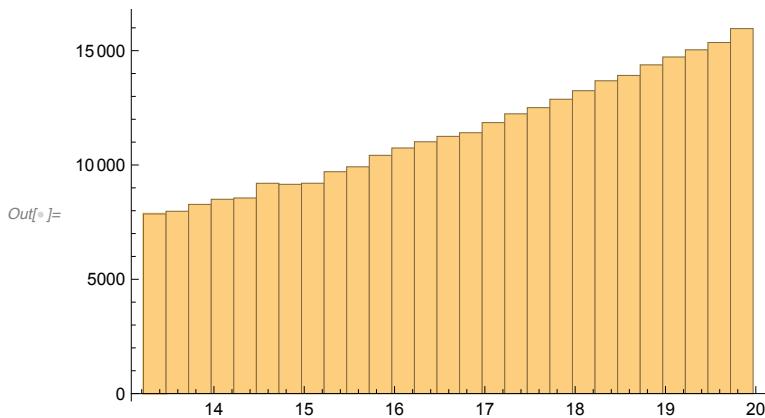
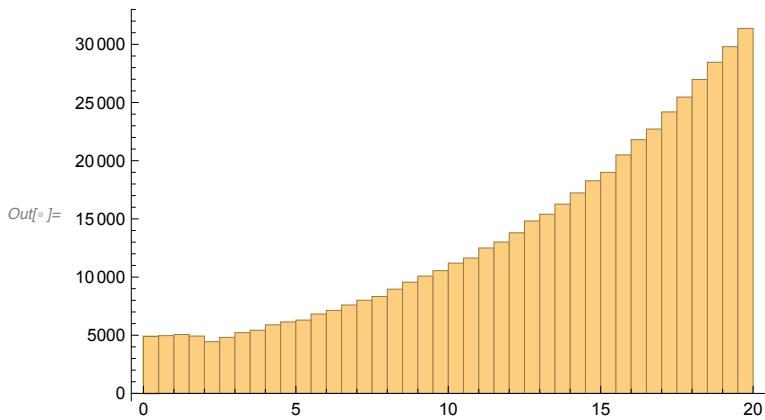
```
l1 = Reap[AvgElectronsBanked[10, 10, 10000]]; l1[[1]]
Histogram[l1[[2, 1]], {0, 10, 0.25}]
Histogram[l1[[2, 1]], {(10 - 2.16401), 10, 0.1}]
```

Out[\circ] = 12.65 ± 0.06



```
l2 = Reap[AvgElectronsBanked[20, 10, 10000]]; l2[[1]]
Histogram[l2[[2, 1]], {0, 20, 0.5}]
Histogram[l2[[2, 1]], {20 - 6.78, 20, 0.25}]
```

Out[=] 7.25 ± 0.04

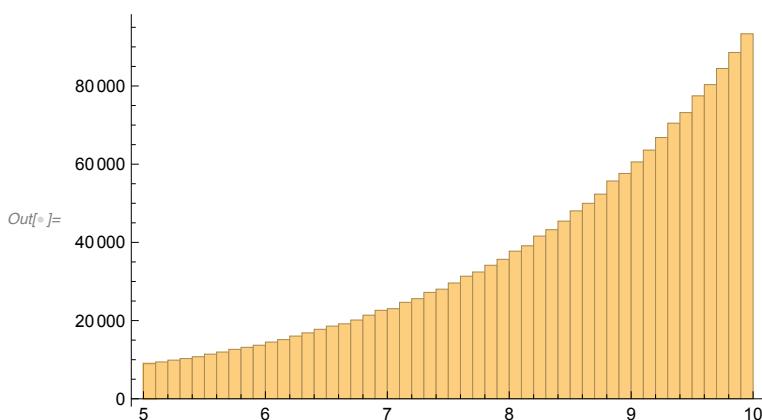
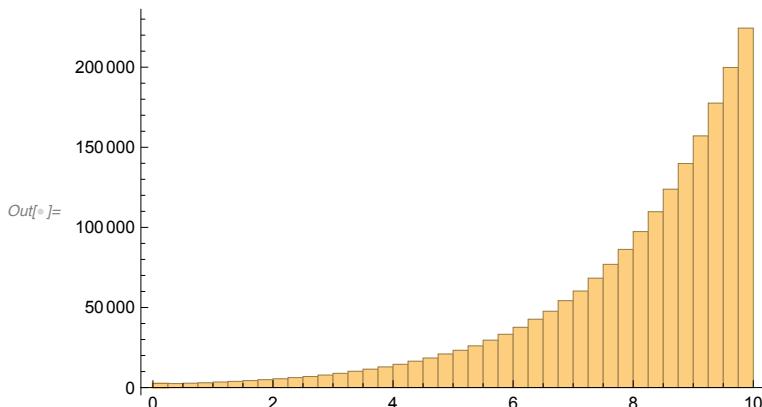


```

l3 = Reap[AvgElectronsBanked[10, 20, 10000]]; l3[[1]]
Histogram[l3[[2, 1]], {0, 10, 0.25}]
Histogram[l3[[2, 1]], {10 - 5, 10, 0.1}]

```

Out[6]= 95.4 ± 0.5



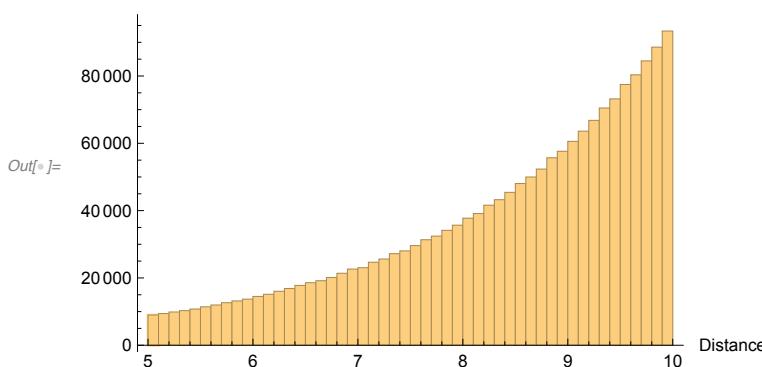
```

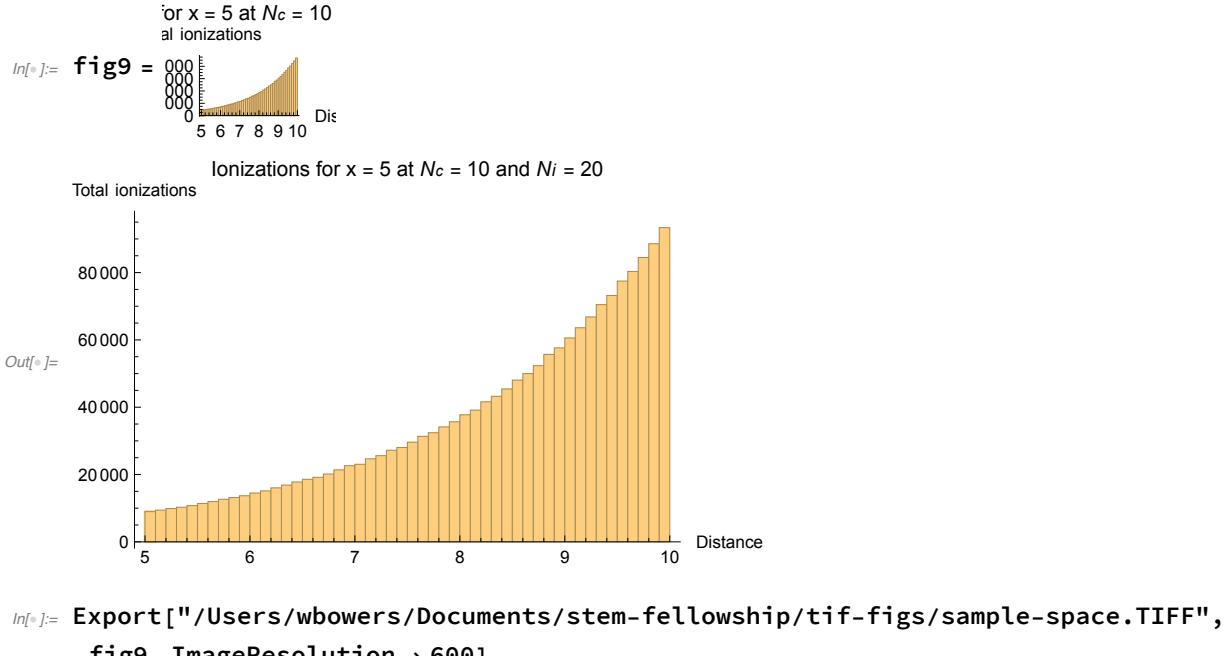
In[7]:= Histogram[l3[[2, 1]], {10 - 5, 10, 0.1}, AxesLabel -> {"Distance", "Total ionizations"}, PlotLabel -> "Ionizations for x = 5 at Nc = 10 and Ni = 20"]

```

Ionizations for x = 5 at $N_c = 10$ and $N_i = 20$

Total ionizations





Out[6]= /Users/wbowers/Documents/stem-fellowship/tif-figs/sample-space.TIFF

Parallel plate breakdown 1D

```

In[7]:= FindBreakdown[r_, \gamma_, reps_] := Module[{alpha, idx, Nc, Ni},
alpha = FindAlpha[r, reps][[1]];
Nc = Log[\gamma] / alpha + r;
Ni = (1/r) (Nc);
{Nc, Ni}]

```

```

In[8]:= bdpp1d = Table[FindBreakdown[r, 50, 500], {r, 1/50, 5, 1/4}]

```

```

Out[8]= {{4.38 \pm 0.24, 219. \pm 12.}, {6.51 \pm 0.29, 24.1 \pm 1.1}, {9.03 \pm 0.11, 17.37 \pm 0.21}, {11.81 \pm 0.10, 15.34 \pm 0.13}, {15.19 \pm 0.09, 14.90 \pm 0.09}, {19.43 \pm 0.20, 15.30 \pm 0.15}, {24.9 \pm 0.4, 16.38 \pm 0.29}, {30.7 \pm 0.5, 17.33 \pm 0.27}, {38.6 \pm 0.7, 19.1 \pm 0.4}, {48.3 \pm 0.6, 21.26 \pm 0.25}, {59.6 \pm 0.8, 23.67 \pm 0.30}, {73.3 \pm 1.7, 26.5 \pm 0.6}, {84.8 \pm 3.5, 28.1 \pm 1.1}, {113. \pm 4., 34.6 \pm 1.3}, {153. \pm 7., 43.4 \pm 2.0}, {185. \pm 6., 49.1 \pm 1.5}, {246. \pm 13., 61.3 \pm 3.2}, {311. \pm 19., 73. \pm 4.}, {406. \pm 38., 90. \pm 8.}, {447. \pm 34., 94. \pm 7.}}

```

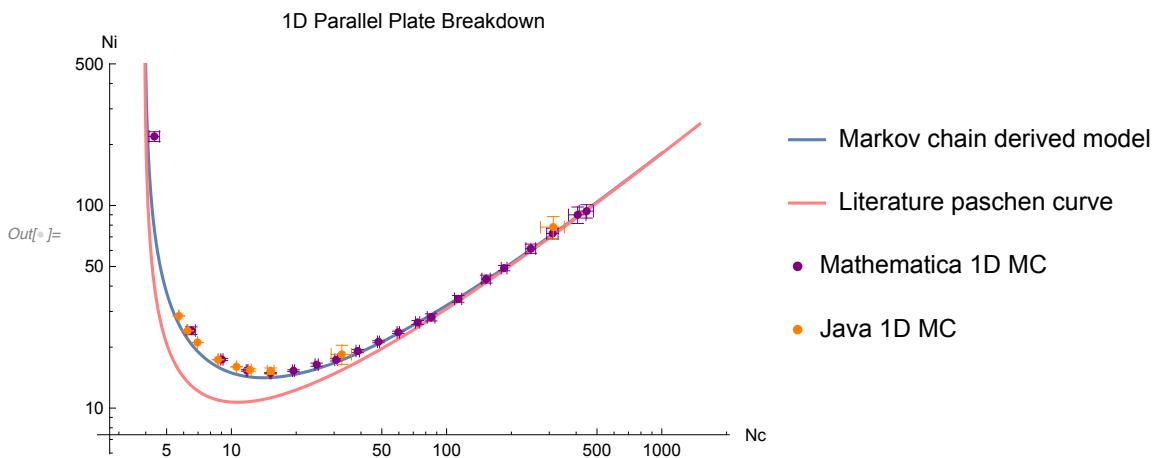
```
In[1]:= javabdd1d = {{Around[5.700231943597744, 0.01207241166528697],  
 Around[28.50115971798872, 0.06036205832643485]},  
 {Around[6.259659518165803, 0.01294585426723786],  
 Around[24.075613531406933, 0.049791747181684075]},  
 {Around[6.956648887715002, 0.006510191038770662],  
 Around[21.08075420519698, 0.01972785163263837]},  
 {Around[8.666186577739973, 0.01778912290701574],  
 Around[17.332373155479946, 0.03557824581403148]},  
 {Around[10.558215754398047, 0.018609784723081196],  
 Around[15.997296597572799, 0.028196643519819993]},  
 {Around[12.189849737655585, 0.1691424999247856],  
 Around[15.430189541336182, 0.2141044302845387]},  
 {Around[15.233034583855726, 0.5468246433362867],  
 Around[15.233034583855726, 0.5468246433362867]},  
 {Around[32.607032745474434, 3.5391212217852397],  
 Around[18.422052398573125, 1.9995035151329037]},  
 {Around[37.56666730405342, 4.788665312298926], Around[19.56597255419449,  
 2.494096516822358]} } {Around[226.1883191276727, 34.15818691413011],  
 Around[61.13197814261424, 9.231942409224354]}, {Around[312.9949279966182,  
 39.66595313492955], Around[78.24873199915454, 9.916488283732388]}}
```

Out[1]= $\left\{ \{5.700 \pm 0.012, 28.50 \pm 0.06\}, \{6.260 \pm 0.013, 24.08 \pm 0.05\}, \{6.957 \pm 0.007, 21.081 \pm 0.020\}, \{8.666 \pm 0.018, 17.33 \pm 0.04\}, \{10.558 \pm 0.019, 15.997 \pm 0.028\}, \{12.19 \pm 0.17, 15.43 \pm 0.21\}, \{15.2 \pm 0.5, 15.2 \pm 0.5\}, \{32.6 \pm 3.5, 18.4 \pm 2.0\}, \{(8.5 \pm 1.7) \times 10^3, 1196. \pm 236.\}, \{313. \pm 40., 78. \pm 10.\} \right\}$

```
In[2]:= Solve[Log[51] == Ni * ProductLog[Nc / Ni * Exp[-Nc / Ni]], Ni]
```

```
Out[2]=  $\left\{ \left\{ Ni \rightarrow \frac{-Nc - \text{Log}[51]}{\text{Log}\left[\frac{\text{Log}[51]}{Nc}\right]} \right\} \right\}$ 
```

```
In[6]:= markovmodelfig = Show[LogLogPlot[-Nc - Log[51], Log[Log[51]/Nc], {Nc, 3, 1000}, PlotLegends -> {"Markov chain derived model"}], LogLogPlot[Nc / (Log[Nc] - Log[Log[51]]), {Nc, 3, 1500}, PlotStyle -> Pink, PlotLegends -> {"Literature paschen curve"}], ListLogLogPlot[bdpp1d, PlotLegends -> {"Mathematica 1D MC"}, PlotStyle -> Purple], ListLogLogPlot[javabd1d, PlotStyle -> Orange, PlotLegends -> {"Java 1D MC"}], AxesLabel -> {"Nc", "Ni"}, PlotLabel -> "1D Parallel Plate Breakdown", PlotRange -> {{1, 7.5}, {2, 6}}, AxesOrigin -> {1, 2}]
```



```
In[7]:= Export["/Users/wbowers/Documents/stem-fellowship/figures/1d-pp-breakdown.TIFF", markovmodelfig, ImageResolution -> 600]
Out[7]= /Users/wbowers/Documents/stem-fellowship/figures/1d-pp-breakdown.TIFF
```

Parallel plate breakdown 3D

```
FindBreakdown3D[d_, γ_, tbound_, Ni0_, reps_] :=
Module[{region, list, i, errors, max, params, fit,
  error, p1, p2, p3, p4, f, vals, errorslist, listForFit},
  list =
  Table[{Ni, AvgElectrons3D[d, Ni, d, 15, tbound, reps]}, {Ni, Ni0 - 3, Ni0 + 3}];
  listForFit = {#[1], #[2, 1]} & /@ list;
  errorslist = #[2, 2] & /@ list;
  fit = NonlinearModelFit[listForFit,
    p1 Exp[p2 x] + p3, {p1, p2, p3}, x, Weights -> 1/errorslist^2];
  f = (Log[(γ - p3)/p1]/p2);
  error = {D[f, p1], D[f, p2], D[f, p3]}.fit["CovarianceMatrix"].
    Transpose[{D[f, p1], D[f, p2], D[f, p3]}] /. fit["BestFitParameters"];
  Around[f, Sqrt[error]] /. fit["BestFitParameters"]
]
```

```

In[1]:= FindBreakdown3Dv2[4.5, 50, 0.8, 30, 3000]
{{27, 45.5 ± 0.7}, {28, 46.1 ± 0.7}, {29, 47.5 ± 0.7},
{30, 48.5 ± 0.7}, {31, 48.7 ± 0.8}, {32, 49.9 ± 0.8}, {33, 51.3 ± 0.8} }

Out[1]= 32.01 ± 0.21

In[2]:= left = Table[{d, FindBreakdown3Dv2[d, 50, 0.8, 20, 500]}, {d, 5, 6, 0.5}]
Out[2]= {{5., 21.2 ± 0.4}, {5.5, 16.75 ± 0.35}, {6., 11. ± 5.} }

In[3]:= min = Table[{d, FindBreakdown3Dv2[d, 50, 0.8, 10, 100]}, {d, 6, 16.5, 2}]
Out[3]= {{6, 14.4 ± 0.4}, {8, 11.73 ± 0.13}, {10, 10.790 ± 0.034},
{12, 10.60 ± 0.04}, {14, 10.52 ± 0.07}, {16, 10.70 ± 0.05} }

In[4]:= right2 = Table[{d, FindBreakdown3Dv2[d, 50, 0.8, 13, 50]}, {d, 17, 57, 4}]
Out[4]= {{17, 10.81 ± 0.08}, {21, 11.24 ± 0.04}, {25, 12.07 ± 0.09},
{29, 12.66 ± 0.05}, {33, 13.45 ± 0.05}, {37, 14.255 ± 0.029}, {41, 14.939 ± 0.018},
{45, 15.66 ± 0.09}, {49, 16.49 ± 0.22}, {53, 17.48 ± 0.09}, {57, 17.74 ± 0.30} }

In[5]:= right3 = Table[{d, FindBreakdown3Dv2[d, 50, 0.8, 17, 25]}, {d, 60, 100, 8}]
Out[5]= {{60, 18.52 ± 0.09}, {68, 19.97 ± 0.07}, {76, 21.31 ± 0.35},
{84, 23.7 ± 1.2}, {92, 24.7 ± 0.9}, {100, 26.2 ± 1.4} }

In[6]:= bdfinal = {{4.5, 32.01 ± 0.21}, {5., 21.2 ± 0.4},
{5.5, 16.75 ± 0.35}, {6, 14.4 ± 0.4}, {8, 11.73 ± 0.13}, {10, 10.790 ± 0.034},
{12, 10.60 ± 0.04}, {14, 10.52 ± 0.07}, {16, 10.70 ± 0.05}, {17, 10.77 ± 0.06},
{20, 11.18 ± 0.06}, {23, 11.64 ± 0.07}, {26, 12.21 ± 0.06}, {32, 13.27 ± 0.05},
{37, 14.255 ± 0.029}, {41, 14.939 ± 0.018}, {49, 16.49 ± 0.22},
{60, 18.52 ± 0.09}, {76, 21.31 ± 0.35}, {92, 24.7 ± 0.9}, {100, 26.2 ± 1.4} }

Out[6]= {{4.5, 32.01 ± 0.21}, {5., 21.2 ± 0.4}, {5.5, 16.75 ± 0.35}, {6, 14.4 ± 0.4}, {8, 11.73 ± 0.13},
{10, 10.790 ± 0.034}, {12, 10.60 ± 0.04}, {14, 10.52 ± 0.07}, {16, 10.70 ± 0.05},
{17, 10.77 ± 0.06}, {20, 11.18 ± 0.06}, {23, 11.64 ± 0.07}, {26, 12.21 ± 0.06},
{32, 13.27 ± 0.05}, {37, 14.255 ± 0.029}, {41, 14.939 ± 0.018}, {49, 16.49 ± 0.22},
{60, 18.52 ± 0.09}, {76, 21.31 ± 0.35}, {92, 24.7 ± 0.9}, {100, 26.2 ± 1.4} }

In[7]:= javabd3d = {{Around[3.9349346617986107`, 0.004098654929587719], 100},
{7., Around[12.5973499575545, 0.021042290566655936`]}, 
{10., Around[10.665900655226146`, 0.01147892717515875`]}, 
{20., Around[11.199362729457336`, 0.0055071173430502485`]}, 
{40., Around[14.831063909226039`, 0.011373607275817859`]}, 
{80., Around[22.424620315119526`, 0.003938560064511855`]}, 
{160., Around[36.60786133347448, 0.009721595486457082]}} 

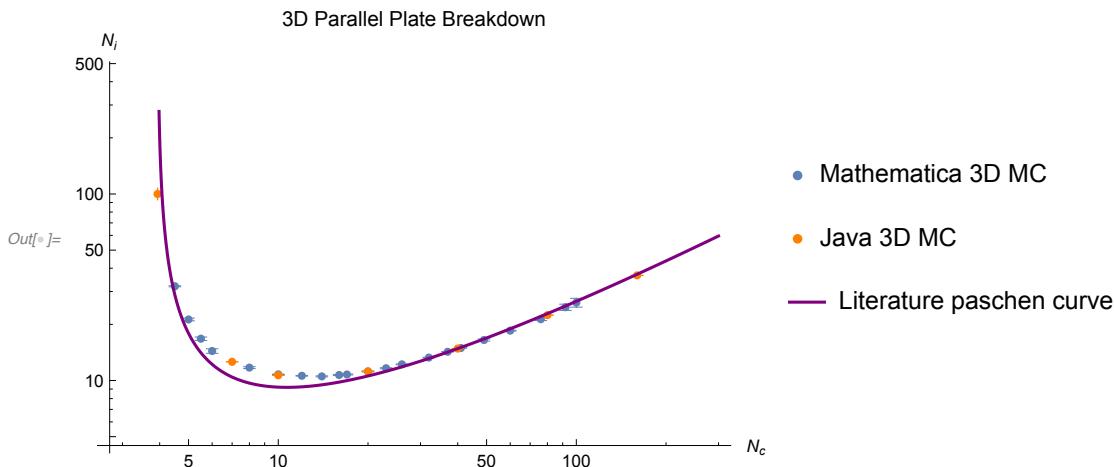
Out[7]= {{3.935 ± 0.004, 100}, {7., 12.597 ± 0.021}, {10., 10.666 ± 0.011}, {20., 11.199 ± 0.006},
{40., 14.831 ± 0.011}, {80., 22.425 ± 0.004}, {160., 36.608 ± 0.010} }

```

```
In[6]:= fig10 = Show[ListLogLogPlot[bdfinal, PlotLegends -> {"Mathematica 3D MC"}],
ListLogLogPlot[javabd3d, PlotStyle -> Orange, PlotLegends -> {"Java 3D MC"}],
LogLogPlot[0.86 * Nc / (Log[Nc] - Log[Log[51]]), {Nc, 3.9, 300}, PlotStyle -> Purple,
PlotLegends -> {"Literature paschen curve"}], PlotRange -> {1.5, 6},
```

```
AxesOrigin -> {1, 1.5}, AxesLabel -> {Subscript[N, c], Subscript[N, i]},
```

```
PlotLabel -> "3D Parallel Plate Breakdown"]
```



```
In[7]:= Export["/Users/wbowers/Documents/stem-fellowship/figures/3d-pp-breakdown.TIFF",
fig10, ImageResolution -> 600]
```

```
Out[7]= /Users/wbowers/Documents/stem-fellowship/figures/3d-pp-breakdown.TIFF
```

Sphere-in-sphere breakdown

```
In[8]:= FindBreakdownSS[Nc_, γ_, Ni0_, dt_, reps_] :=
Module[{region, list, i, errors, max, params, fit,
error, p1, p2, p3, p4, f, vals, errorslist, listForFit},
list = Table[{Ni, AvgElectronsEuler[Nc, Ni, 5, 15, dt, 4/5, False, reps]},
{Ni, Ni0 - 3, Ni0 + 2}];
Print[list];
listForFit = {#[[1]], #[[2, 1]]} & /@ list;
errorslist = #[[2, 2]] & /@ list;
fit = NonlinearModelFit[listForFit,
p1 Exp[p2 x] + p3, {p1, p2, p3}, x, Weights -> 1 / errorslist^2];
f = (Log[(γ - p3) / p1] / p2);
error = {D[f, p1], D[f, p2], D[f, p3]}.fit["CovarianceMatrix"].
Transpose[{D[f, p1], D[f, p2], D[f, p3]}] /. fit["BestFitParameters"];
Around[f, Sqrt[error]] /. fit["BestFitParameters"]
]
```

```

In[1]:= FindNC[Ni_, γ_, Nc0_, dt_, reps_] :=
  Module[{region, list, i, errors, max, params, fit, error,
    p1, p2, p3, p4, f, vals, errorslist, listForFit},
    list = Table[{Nc, AvgElectronsEuler[Nc, Ni, 5, 15, dt, 4/5, False, reps]}, {
      Nc, Nc0 - 1, Nc0 + 1, 1/2}];
    Print[list];
    listForFit = {#[[1]], #[[2, 1]]} & /@ list;
    errorslist = #[[2, 2]] & /@ list;
    fit = NonlinearModelFit[listForFit,
      p1 Exp[p2 x] + p3, {p1, p2, p3}, x, Weights → 1/errorslist^2];
    f = (Log[(γ - p3) / p1] / p2);
    error = {D[f, p1], D[f, p2], D[f, p3]}.fit["CovarianceMatrix"];
    Transpose[{D[f, p1], D[f, p2], D[f, p3]}] /. fit["BestFitParameters"];
    Around[f, Sqrt[error]] /. fit["BestFitParameters"]
  ]
]

In[2]:= FindNC[100, 50, 4, 1/5, 300]
{ {3, 18.9 ± 1.2}, {7/2, 34.4 ± 1.9}, {4, 50.1 ± 2.8}, {9/2, 77. ± 4.}, {5, 130. ± 7.} }

Out[2]= 3.97 ± 0.05

In[3]:= FindBreakdownSS[5, 50, 25, 2/5, 300]
{ {22, 47.0 ± 2.2}, {23, 44.0 ± 2.2}, {24, 45.9 ± 2.5},
  {25, 50.3 ± 2.3}, {26, 53.8 ± 2.6}, {27, 50.6 ± 2.8}, {28, 58.9 ± 3.2} }

Out[3]= 25.6 ± 0.8

In[4]:= FindBreakdownSS[7, 50, 15, 2/5, 50]
{ {12, 38.2 ± 3.4}, {13, 45. ± 4.}, {14, 53. ± 6.},
  {15, 58. ± 6.}, {16, 81. ± 8.}, {17, 73. ± 9.}, {18, 93. ± 8.} }

Out[4]= 13.58 ± 0.31

In[5]:= FindBreakdownSS[12, 50, 8, 2/5, 50]
{ {5, 4.92 ± 0.18}, {6, 7.76 ± 0.29}, {7, 12.6 ± 0.5}, {8, 18.2 ± 0.7}, {9, 25.8 ± 1.2}, {10, 40.9 ± 2.5} }

Out[5]= 10.65 ± 0.15

In[6]:= FindBreakdownSS[30, 50, 13, 2/5, 30]
{ {10, 16.9 ± 1.1}, {11, 27.4 ± 1.5}, {12, 43.0 ± 2.8}, {13, 66. ± 4.}, {14, 99. ± 6.}, {15, 160. ± 9.} }

Out[6]= 12.36 ± 0.04

In[7]:= FindBreakdownSS[50, 50, 15, 2/5, 20]
{ {12, 13.4 ± 0.9}, {13, 21.9 ± 1.6}, {14, 31.3 ± 2.2}, {15, 59. ± 4.}, {16, 87. ± 5.}, {17, 129. ± 11.} }

Out[7]= 14.83 ± 0.09

```

```
In[1]:= FindBreakdownSS[90, 50, 19, 2/5, 20]
{{16, 13.4 ± 1.4}, {17, 21.4 ± 1.8}, {18, 36.9 ± 2.3}, {19, 45.6 ± 3.0}, {20, 71. ± 5.}, {21, 110. ± 8.}}
Out[1]= 19.03 ± 0.12

In[2]:= FindBreakdownSS[150, 50, 24, 2/5, 10]
{{21, 14.8 ± 2.5}, {22, 20.5 ± 1.5}, {23, 23. ± 5.}, {24, 37. ± 5.}, {25, 42. ± 7.}, {26, 81. ± 11.}}
Out[2]= 25.00 ± 0.20

In[3]:= ssbd = {{4.036 ± 0.018, 100}, {5, 25.6 ± 0.8}, {7, 13.58 ± 0.31}, {12, 10.65 ± 0.15},
{30, 12.36 ± 0.04}, {50, 14.83 ± 0.09}, {90, 19.03 ± 0.12}, {150, 25.00 ± 0.20}}
Out[3]= {{4.036 ± 0.018, 100}, {5, 25.6 ± 0.8}, {7, 13.58 ± 0.31}, {12, 10.65 ± 0.15},
{30, 12.36 ± 0.04}, {50, 14.83 ± 0.09}, {90, 19.03 ± 0.12}, {150, 25.00 ± 0.20}}

In[4]:= drwhitmerdata =
{{6.494 ± 0.019, 15.}, {5.469 ± 0.017, 20.}, {10, 11.227 ± 0.008}, {17.8, 11.026 ± 0.009},
{31.6, 12.636 ± 0.006}, {56.2, 15.558 ± 0.020}, {100., 20.350 ± 0.025}}
Out[4]= {{6.494 ± 0.019, 15.}, {5.469 ± 0.017, 20.}, {10, 11.227 ± 0.008}, {17.8, 11.026 ± 0.009},
{31.6, 12.636 ± 0.006}, {56.2, 15.558 ± 0.020}, {100., 20.350 ± 0.025}]

In[5]:= fig11 = Show[ListLogLogPlot[ssbd, PlotLegends -> {"Mathematica MC"}],
ListLogLogPlot[drwhitmerdata, PlotStyle -> Orange, PlotLegends -> {"C++ MC"}],
LogLogPlot[Nc / (Log[Nc] - Log[Log[51]]), {Nc, 0, 300}, PlotStyle -> Purple,
PlotLegends -> {"Literature parallel plate paschen curve"}], AxesLabel -> {Nc, Ni},
PlotLabel -> "Sphere-in-sphere Paschen Curve", PlotRange -> All]
Sphere-in-sphere Paschen Curve
Ni
100
50
10
Out[5]=
5
10
50
100
Nc
• Mathematica MC
● C++ MC
— Literature parallel plate paschen curve
```

```
In[6]:= Export["/Users/wbowers/Documents/High-res-fusor-figures/ss-breakdown.png",
fig11, ImageResolution -> 300]
Out[6]= /Users/wbowers/Documents/High-res-fusor-figures/ss-breakdown.png
```

Sphere-in-cylinder breakdown

```
In[1]:= FindBreakdownNISC[\lambda_, \gamma_, Ni0_, reps_] :=
Module[{region, list, i, errors, max, params, fit,
  error, p1, p2, p3, p4, f, vals, errorslist, listForFit},
  list =
  ParallelTable[{Ni, MaxElectrons[\lambda, Ni, reps]}, {Ni, Ni0 - 3/2, Ni0 + 3/2, 1/2}];
  Print[list];
  listForFit = {#[[1]], #[[2, 1]]} & /@ list;
  errorslist = #[[2, 2]] & /@ list;
  fit = NonlinearModelFit[listForFit,
    p1 Exp[p2 x] + p3, {p1, p2, p3}, x, Weights \rightarrow 1/errorslist^2];
  f = (Log[(\gamma - p3) / p1] / p2);
  error = {D[f, p1], D[f, p2], D[f, p3]}.fit["CovarianceMatrix"].
    Transpose[{D[f, p1], D[f, p2], D[f, p3]}] /. fit["BestFitParameters"];
  Around[f, Sqrt[error]] /. fit["BestFitParameters"]
]

In[2]:= FindBreakdownNCSC[Ni_, \gamma_, Nc0_, reps_] :=
Module[{region, list, i, errors, max, params, fit,
  error, p1, p2, p3, p4, f, vals, errorslist, listForFit},
  list = ParallelTable[
    {Nc, MaxElectrons[53 / (10 * Nc), Ni, reps]}, {Nc, Nc0 - 1, Nc0 + 1, 1/2}];
  Print[list];
  listForFit = {#[[1]], #[[2, 1]]} & /@ list;
  errorslist = #[[2, 2]] & /@ list;
  fit = NonlinearModelFit[listForFit,
    p1 Exp[p2 x] + p3, {p1, p2, p3}, x, Weights \rightarrow 1/errorslist^2];
  f = (Log[(\gamma - p3) / p1] / p2);
  error = {D[f, p1], D[f, p2], D[f, p3]}.fit["CovarianceMatrix"].
    Transpose[{D[f, p1], D[f, p2], D[f, p3]}] /. fit["BestFitParameters"];
  Around[f, Sqrt[error]] /. fit["BestFitParameters"]
]

In[3]:= FindBreakdownNCSC[40, 50, 34/10, 500]
Out[3]= 0.63622
Out[4]= 0.645566
Out[5]= -0.746137
Out[6]= -0.756372
Out[7]= -0.665681
Out[8]= {12/5, 17.6 \pm 1.3}, {29/10, 31.3 \pm 2.1}, {17/5, 47.4 \pm 3.4}, {39/10, 83. \pm 6.}, {22/5, 101. \pm 7.}
Out[9]= 3.41 \pm 0.07
```

NC = 10

In[]:= **FindBreakdownNCSC[30, 50, 4, 500]**

(kernel 4) **0.659182**

(kernel 3) **0.754227**

(kernel 4) **-0.645615**

(kernel 2) **0.691999**

(kernel 1) **-0.708711**

During evaluation of *In[1047]:=*

$$\left\{ \{3, 26.9 \pm 1.9\}, \left\{ \frac{7}{2}, 41.9 \pm 2.7 \right\}, \{4, 74. \pm 5.\}, \left\{ \frac{9}{2}, 92. \pm 6. \right\}, \{5, 128. \pm 9.\} \right\}$$

Out[]= **3.63 ± 0.07**

In[]:= **FindBreakdownNCSC[18, 50, 4, 500]**

(kernel 4) **-0.858654**

(kernel 3) **-0.599666**

(kernel 4) **0.731309**

(kernel 2) **-0.75725**

(kernel 1) **-0.616007**

During evaluation of *In[1043]:=*

$$\left\{ \{3, 17.8 \pm 1.4\}, \left\{ \frac{7}{2}, 30.2 \pm 2.1 \right\}, \{4, 31.0 \pm 2.1\}, \left\{ \frac{9}{2}, 52. \pm 4. \right\}, \{5, 60. \pm 4.\} \right\}$$

Out[]= **4.62 ± 0.18**

In[]:= **FindBreakdownNISC[53 / 100, 50, 11.1, 1000]**

(kernel 4) **-1.**

(kernel 3) **0.367845**

(kernel 2) **-6.29888 × 10⁻⁹**

(kernel 1) **-1.**

(kernel 4) **0.938665**

(kernel 3) **0.22094**

(kernel 2) **0.936544**

During evaluation of *In[978]:=*

$$\{\{9.6, 35.1 \pm 3.1\}, \{10.1, 41.1 \pm 1.6\}, \{10.6, 43.8 \pm 1.2\}, \\ \{11.1, 50.7 \pm 1.3\}, \{11.6, 58.2 \pm 2.3\}, \{12.1, 66.6 \pm 3.3\}, \{12.6, 99. \pm 10.\}\}$$

Out[]= **11.09 ± 0.07**

```

In[6]:= electronvals = {{9.6` , 35.1 ± 3.1}, {10.1` , 41.1 ± 1.6},
{10.6` , 43.8 ± 1.2}, {11.1` , 50.7 ± 1.3}, {11.6` , 58.2 ± 2.3}, {12.1` , 66.6 ± 3.3}}
listForFit = #[[1]], #[[2, 1]] & /@ electronvals;
errorslist = #[[2, 2]] & /@ electronvals;
fit = NonlinearModelFit[listForFit,
p1 Exp[p2 x] + p3, {p1, p2, p3}, x, Weights → 1 / errorslist^2];
f = (Log[(50 - p3) / p1] / p2);
error = {D[f, p1], D[f, p2], D[f, p3]}.fit["CovarianceMatrix"].
Transpose[{D[f, p1], D[f, p2], D[f, p3]}] /. fit["BestFitParameters"];
res = Around[f, Sqrt[error]] /. fit["BestFitParameters"]

Out[6]= {{9.6, 35.1 ± 3.1}, {10.1, 41.1 ± 1.6}, {10.6, 43.8 ± 1.2},
{11.1, 50.7 ± 1.3}, {11.6, 58.2 ± 2.3}, {12.1, 66.6 ± 3.3} }

Out[7]= 11.07 ± 0.04

In[8]:= res[[1]]
Out[8]= 11.0654

In[9]:= fig7 = Show[ListPlot[electronvals], Plot[fit[x], {x, 9, 16}],
Graphics[{Red, PointSize[0.03], Point[{res[[1]], 50}]}], Graphics[{Dashed,
Black, Line[{{9.6, 50}, {res[[1]], 50}}], Line[{{res[[1]], 0}, {res[[1]], 50}}]}],
Graphics[Text["Breakdown Point", {11, 57}]],
AxesLabel → {Subscript["N", "i"], "Electrons"}, PlotLabel → "Finding breakdown for sphere-in-cylinder MC with γ = 50"]

```

Finding breakdown for sphere-in-cylinder MC with $\gamma = 50$

Electrons

Out[9]=

```

In[10]:= Export["/Users/wbowers/Documents/stem-fellowship/figures/scanning-ni.TIFF",
fig7, ImageResolution → 600]

```

Out[10]= /Users/wbowers/Documents/stem-fellowship/figures/scanning-ni.TIFF

```

In[11]:= FindBreakdownNISC[53 / 178, 50, 11.1, 500]

```

```
(kernel 4) -4.36402 × 10-9
(kernel 3) -1.28008 × 10-8
(kernel 2) -1.
(kernel 4) -7.82048 × 10-9
(kernel 1) 2.20918 × 10-33
(kernel 1) General::munfl : 2.20918 × 10-3310 is too small to
represent as a normalized machine number; precision may be lost.
(kernel 1) General::munfl : 2.20918 × 10-3310 is too small to
represent as a normalized machine number; precision may be lost.
(kernel 1) General::munfl : 2.20918 × 10-3310 is too small to
represent as a normalized machine number; precision may be lost.
(kernel 1) General::stop :
Further output of General::munfl will be suppressed during this calculation.
(kernel 3) 1.
(kernel 2) -1.30523 × 10-8
{{9.6, 26.8 ± 0.8}, {10.1, 34.9 ± 1.1}, {10.6, 42.1 ± 1.4}, {11.1, 55. ± 6.}, {11.6, 75. ± 8.}, {12.1, 73.2 ± 2.9}, {12.6, 95. ± 4.}}
Out[=]= 10.99 ± 0.07

In[=]:= FindBreakdownNISC[53 / 200, 50, 13, 100]
Out[=]= $Aborted

(kernel 4) -0.934342
(kernel 3) -5.14173 × 10-9
(kernel 2) -1.
(kernel 4) 1.
(kernel 1) 0.
(kernel 3) -0.912496
(kernel 2) -1.

During evaluation of In[982]:= {{23/2, 55. ± 5.}, {12, 65. ± 11.}, {25/2, 101. ± 9.},
{13, 103. ± 12.}, {27/2, 154. ± 28.}, {14, 183. ± 67.}, {29/2, 180. ± 15.}}
Out[=]= 11.38 ± 0.14

In[=]:= FindBreakdownNISC[53 / 400, 50, 13, 200]
```

```
(kernel 4) -1.
(kernel 3) -4.96763 × 10-9
(kernel 2) 1.
(kernel 4) -0.26817
(kernel 1) -4.1359 × 10-24
(kernel 3) -6.18105 × 10-9
(kernel 2) 1.
```

During evaluation of In[989]:=

$$\left\{ \left\{ \frac{23}{2}, 18.7 \pm 2.6 \right\}, \{12, 23.5 \pm 0.8\}, \left\{ \frac{25}{2}, 29.4 \pm 1.3 \right\}, \{13, 38.4 \pm 2.0\}, \left\{ \frac{27}{2}, 46. \pm 5. \right\}, \{14, 59. \pm 7.\}, \left\{ \frac{29}{2}, 69.0 \pm 3.2 \right\} \right\}$$

Out[989]= 13.680 ± 0.029

In[989]:= **FindBreakdownNISC[53 / 700, 50, 17, 100]**

```
(kernel 4) 0.589819
(kernel 3) -1.
(kernel 4) 0.514641
(kernel 2) 0.214126
(kernel 3) -0.907422
(kernel 1) -0.913193
(kernel 2) 1.
```

During evaluation of In[986]:=

$$\left\{ \left\{ \frac{31}{2}, 28.9 \pm 2.3 \right\}, \{16, 30.6 \pm 2.2\}, \left\{ \frac{33}{2}, 45. \pm 12. \right\}, \{17, 46. \pm 4.\}, \left\{ \frac{35}{2}, 60.9 \pm 2.8 \right\}, \{18, 72. \pm 14.\}, \left\{ \frac{37}{2}, 96. \pm 7. \right\} \right\}$$

Out[986]= 17.12 ± 0.06

In[986]:= **FindBreakdownNISC[53 / 1100, 50, 23, 100]**

```
(kernel 4) -1.
(kernel 3) -0.970182
(kernel 4) -0.306198
(kernel 2) -0.918607
(kernel 3) -0.265945
(kernel 1) -1.03398 × 10-25
(kernel 2) -4.71216 × 10-9
 $\left\{ \left\{ \frac{43}{2}, 56. \pm 15. \right\}, \{22, 60. \pm 4.\}, \left\{ \frac{45}{2}, 87. \pm 7. \right\}, \{23, 100. \pm 6.\}, \left\{ \frac{47}{2}, 124. \pm 11. \right\}, \{24, 163. \pm 9.\}, \left\{ \frac{49}{2}, 214. \pm 16. \right\} \right\}$ 
```

Out[6]= 21.53 ± 0.19

```
In[7]:= FindBreakdownNISC[53 / 1500, 50, 24, 80]
(kernel 4) 2.06795 × 10-25
(kernel 3) -6.58048 × 10-9
(kernel 4) -0.8947
(kernel 2) 0.305773
(kernel 3) -0.355367
(kernel 1) -0.882098
(kernel 2) 0.488683
 $\left\{ \left\{ \frac{45}{2}, 24.3 \pm 2.9 \right\}, \{23, 35. \pm 4.\}, \left\{ \frac{47}{2}, 32.8 \pm 3.2 \right\}, \{24, 41.1 \pm 3.1\}, \left\{ \frac{49}{2}, 48. \pm 5. \right\}, \{25, 54. \pm 5.\}, \left\{ \frac{51}{2}, 61. \pm 7. \right\} \right\}$ 
```

Out[8]= 24.71 ± 0.12

```
In[6]:= scbreakdown = {{5.3 / 3.41 ± 0.07, 40}, {5.3 / 3.63 ± 0.07, 30}, {5.3 / 4.62 ± 0.18, 18}, {53 / 100, 11.09 ± 0.07}, {53 / 200, 11.38 ± 0.14}, {53 / 400, 13.680 ± 0.029}, {53 / 700, 17.12 ± 0.06}, {53 / 1100, 21.53 ± 0.19}, {53 / 1500, 24.71 ± 0.12}}
```

```
sbreakdownscaled = {1 / #[[1]], 15 * #[[2]]} & /@ scbreakdown
```

```
Out[6]= {{1.557 ± 0.031, 40}, {1.459 ± 0.028, 30}, {1.15 ± 0.04, 18}, {53 / 100, 11.09 ± 0.07}, {53 / 200, 11.38 ± 0.14}, {53 / 400, 13.680 ± 0.029}, {53 / 700, 17.12 ± 0.06}, {53 / 1100, 21.53 ± 0.19}, {53 / 1500, 24.71 ± 0.12}}
```

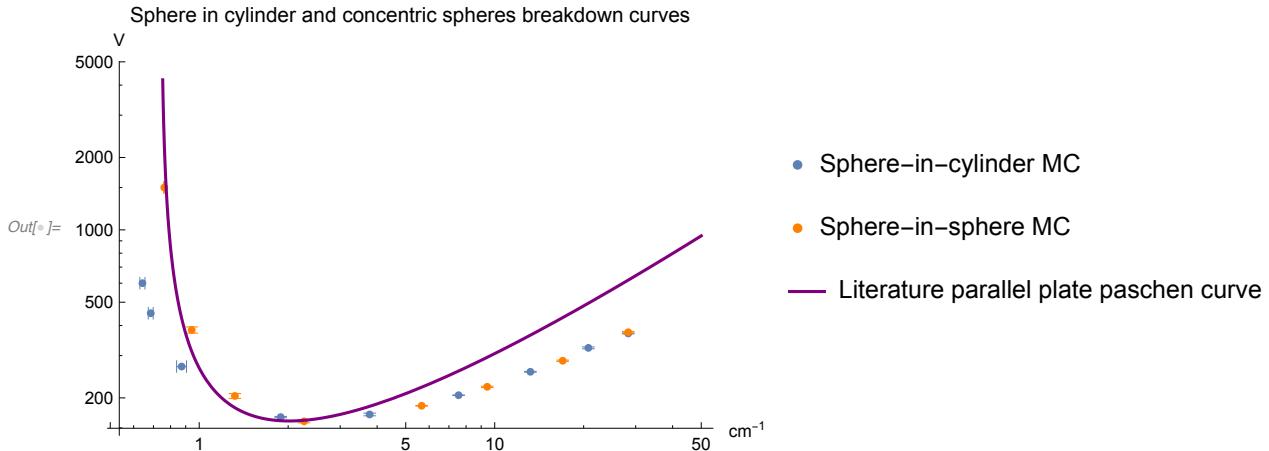
```
Out[7]= {{0.642 ± 0.013, 600}, {0.685 ± 0.013, 450}, {0.872 ± 0.034, 270}, {100 / 53, 166.4 ± 1.0}, {200 / 53, 170.6 ± 2.2}, {400 / 53, 205.2 ± 0.4}, {700 / 53, 256.8 ± 1.0}, {1100 / 53, 322.9 ± 2.8}, {1500 / 53, 370.7 ± 1.7}}
```

```
In[8]:= ssbreakdownscaled = {#[[1]] / 5.3, 15 * #[[2]]} & /@ ssbd
```

```
Out[8]= {{0.7616 ± 0.0035, 1500}, {0.943396, 383. ± 11.}, {1.32075, 204. ± 5.}, {2.26415, 159.8 ± 2.2}, {5.66038, 185.4 ± 0.5}, {9.43396, 222.4 ± 1.4}, {16.9811, 285.4 ± 1.8}, {28.3019, 375.0 ± 3.1}}
```

```
In[9]:= fig12 =
```

```
Show[ListLogLogPlot[sbreakdownscaled, PlotLegends → {"Sphere-in-cylinder MC"}], ListLogLogPlot[ssbreakdownscaled, PlotStyle → Orange, PlotLegends → {"Sphere-in-sphere MC"}], LogLogPlot[15 (5.3) Nc / (Log[Nc * 5.3] - Log[Log[51]]), {Nc, 0, 50}, PlotStyle → Purple, PlotLegends → {"Literature parallel plate paschen curve"}], PlotRange → All, PlotLabel → "Sphere in cylinder and concentric spheres breakdown curves", AxesLabel → {Superscript["cm", "-1"], "V"}]
```



```
In[1]:= Export[
  "/Users/wbowers/Documents/stem-fellowship/tif-figs/ss-and-sc-breakdown.TIFF",
  fig12, ImageResolution → 600]

Out[1]= /Users/wbowers/Documents/stem-fellowship/tif-figs/ss-and-sc-breakdown.TIFF
```

Plotting fusor data

```
In[2]:= data = Drop[Import["/Users/wbowers/Downloads/RAWpashendata.txt", "Table"], 1]
```

```
{ {17.4373, 2086.77}, {17.4627, 2052.92}, {17.4878, 2069.5},
{17.5153, 2036.34}, {17.549, 2081.93}, {17.5836, 2041.17}, {17.6141, 2037.03},
{17.641, 2048.08}, {17.6684, 2031.5}, {17.7051, 2051.54}, {17.7484, 1998.34},
{17.7916, 2043.25}, {17.8291, 2030.12}, {17.8664, 1991.44},
{17.9028, 1962.42}, {17.9676, 1967.26}, ..., {16656.5, 2019.76},
{16660, 2057.06}, {16660, 2035.65}, {16660, 2054.99}, {16660, 2073.64},
{16660, 2073.64}, {16660, 2073.64}, {16663.4, 2073.64}, {16666.9, 2081.24},
{16670.7, 2081.24}, {16677.3, 2079.17}, {16684, 2102.66}, {16690.9, 2073.64},
{16697.9, 2077.79}, {16704.9, 2082.62}, {-----, -----} }
```

large output

show less

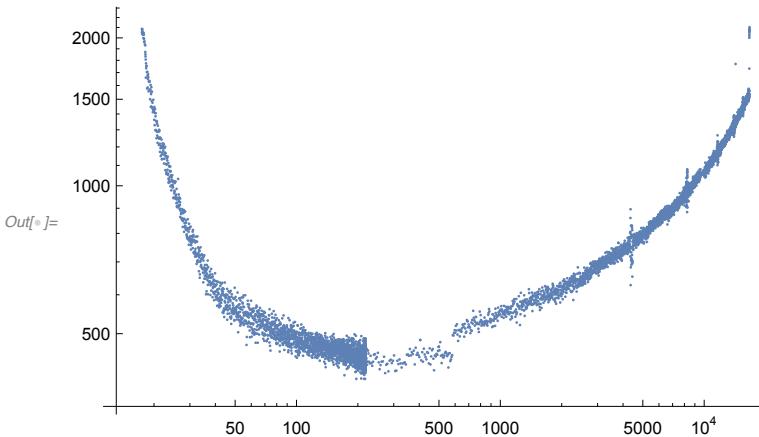
show more

show all

set si

ze limit...

```
In[3]:= ListLogLogPlot[data]
```



```
In[4]:= b = 365
```

```
Out[4]= 365
```

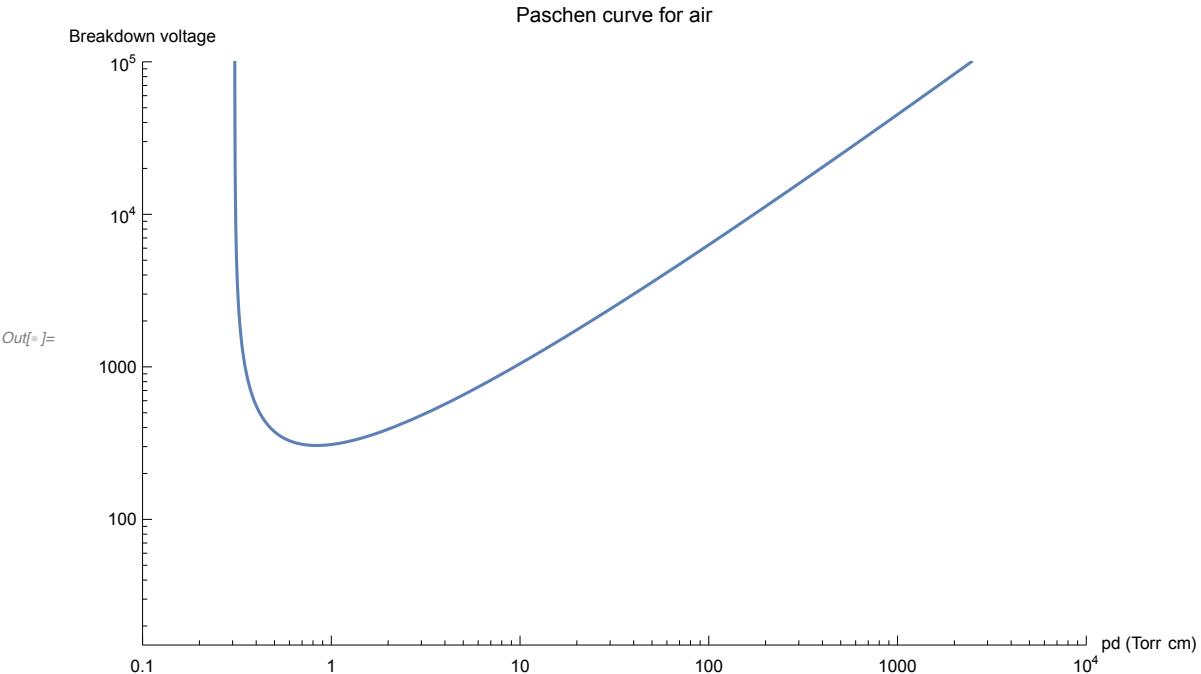
```
In[5]:= a = 15
```

```
Out[5]= 15
```

```
In[6]:= k = Log[a / Log[1 + 1 / (0.01)]]
```

```
Out[6]= 1.17871
```

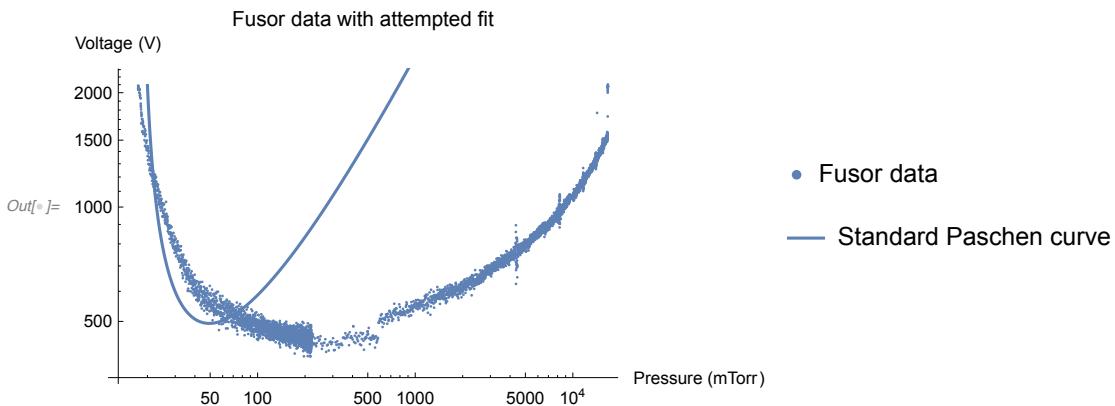
```
In[6]:= fig1 = LogLogPlot[b x / (Log[x] + k),
{x, 0.1, 10000}, PlotRange -> {{0.1, 10000}, {15, 100000}},
AxesOrigin -> {0.1, 15}, PlotLabel -> "Paschen curve for air",
AxesLabel -> {"pd (Torr cm)", "Breakdown voltage"}]
```



```
In[7]:= Export["/Users/wbowers/Documents/stem-fellowship/tif-figs/air-paschen-curve.TIFF",
fig1, ImageResolution -> 600]
```

```
Out[7]= /Users/wbowers/Documents/stem-fellowship/tif-figs/air-paschen-curve.TIFF
```

```
In[8]:= fig2 = Show[ListLogLogPlot[data, PlotLegends -> {"Fusor data"},
PlotStyle -> {PointSize[0.004]}], LogLogPlot[10 x / (Log[x] - 2.9),
{x, 20, 1000}, PlotLegends -> {"Standard Paschen curve"}],
AxesLabel -> {"Pressure (mTorr)", "Voltage (V)"},
PlotLabel -> "Fusor data with attempted fit"]
```



```
In[]:= Export["/Users/wbowers/Documents/stem-fellowship/tif-figs/attempted-fit.TIFF",
  fig2, ImageResolution → 600]

Out[]= /Users/wbowers/Documents/stem-fellowship/tif-figs/attempted-fit.TIFF
```