

Brute Force Simulation

Simulation Configuration

Classical field parameters

636

Number of emitting atoms

293

Temperature (K)

Time resolution (ps)

Maximum time (ns)

f_0 (GHz) = 456811.0

Δf (GHz) = 1.0

σ (GHz) = 7.58

Photon count parameters

Expected counts in time t_{\max}

Total photon counts overall (approximate)

Number of classical simulations: 100

Plotting options

Available Plots:

- Single instance of classical calculations ☒
- Average of all classical calculations ☒
- Total counts from all time windows ☒
- Autocorrelation of photon count time series ☒

Make Plots [☒

Plots

Window:

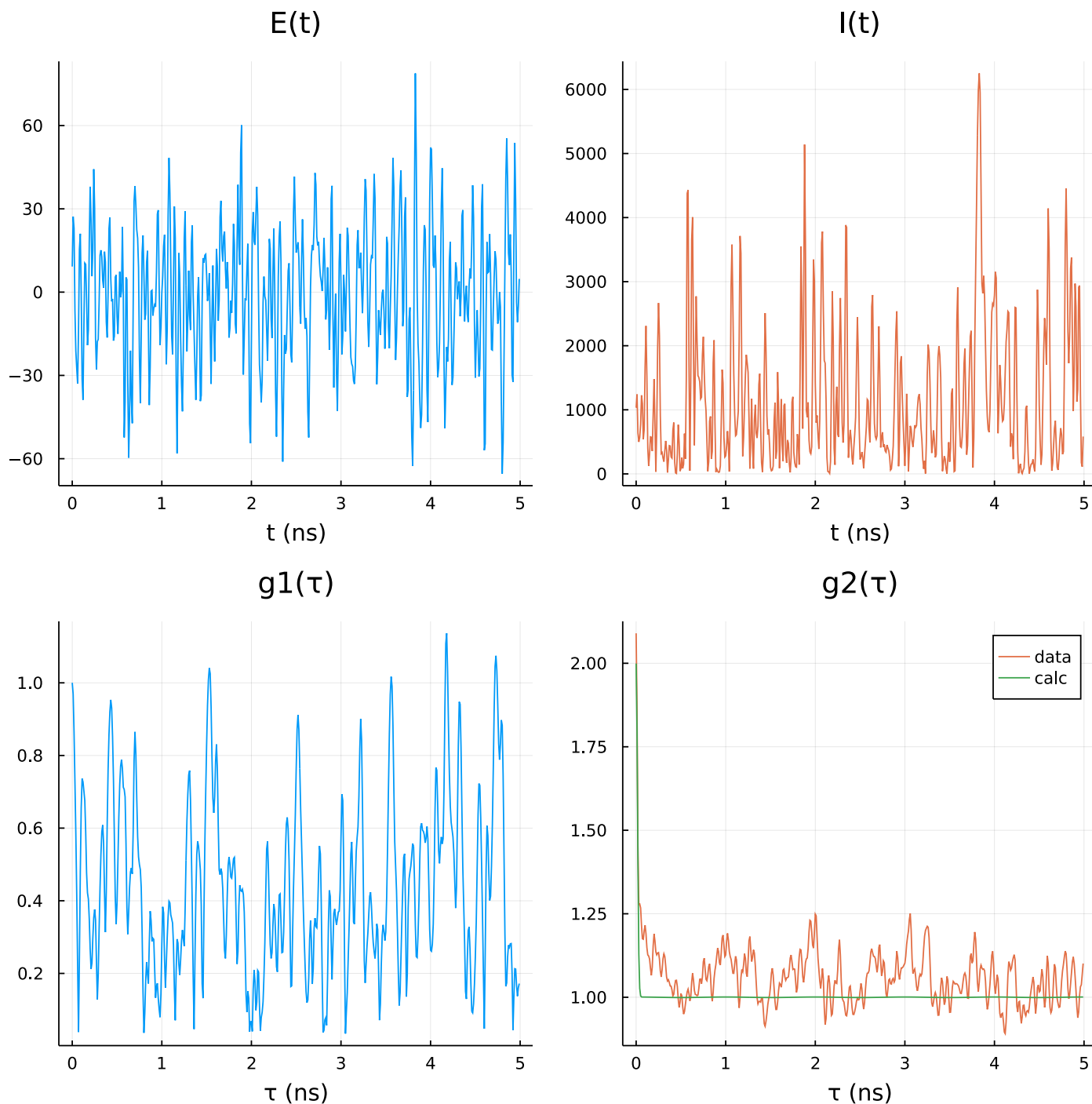
Window size (ns) = 5.0

Window position (ns) = 0.0

Size:

Position:

Single instance of classical calculations:



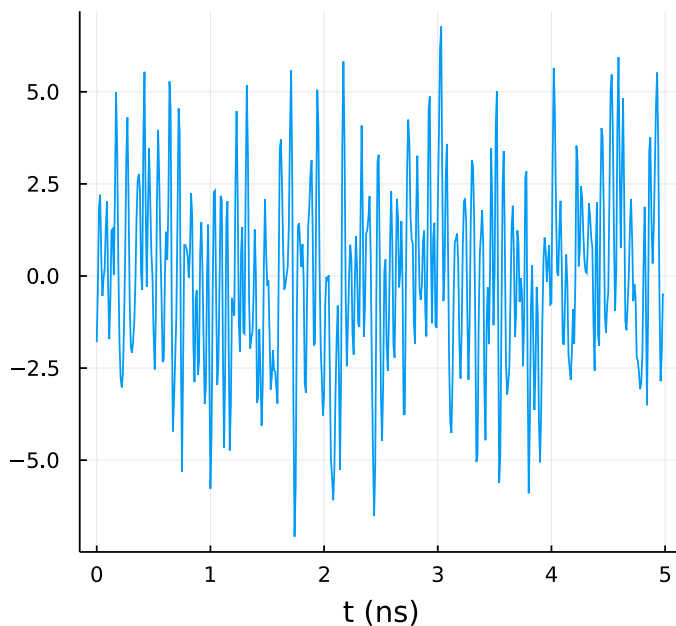
Download data for classical field and correlation plots:

E-field/Intensity: classical_field_single.csv

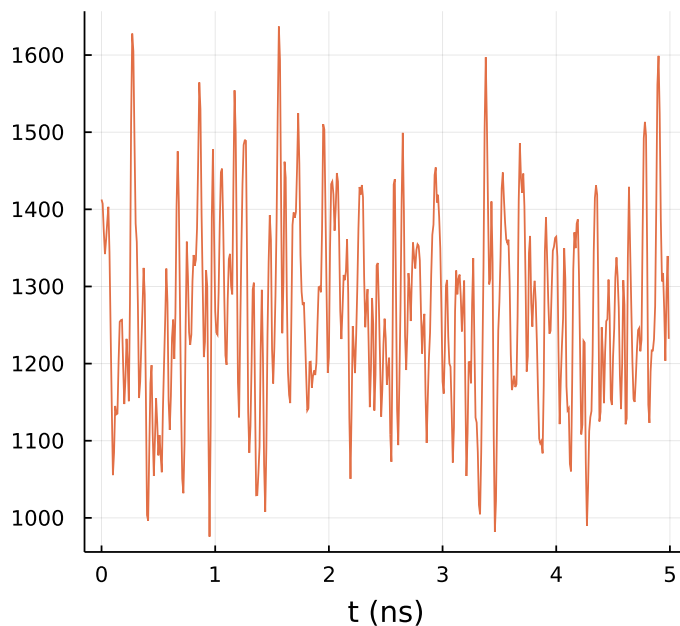
Correlations: classical_correlations_single.csv

Average of all classical calculations:

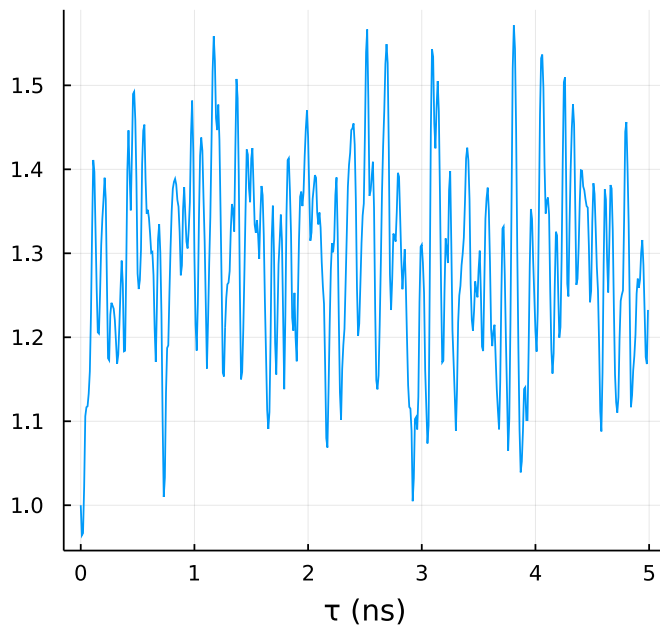
Average $E(t)$: $N = 100$



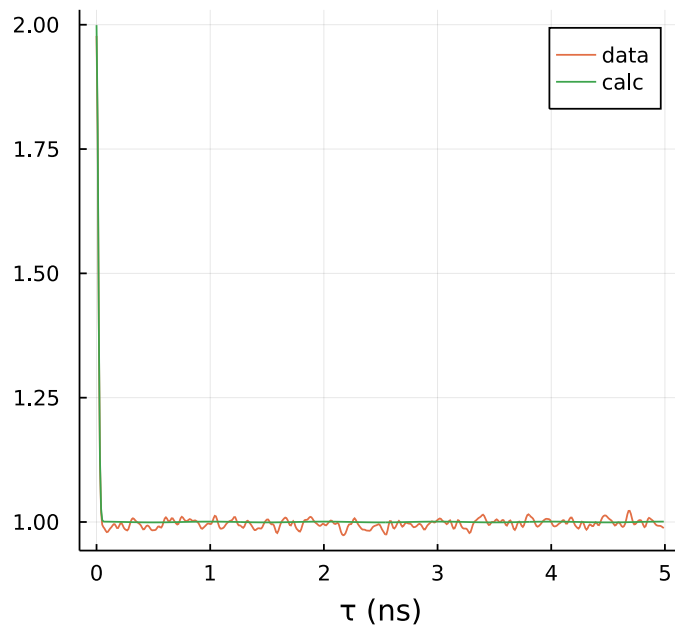
Average $I(t)$: $N = 100$



Average $g_1(\tau)$: $N = 100$



Average $g_2(\tau)$: $N = 100$

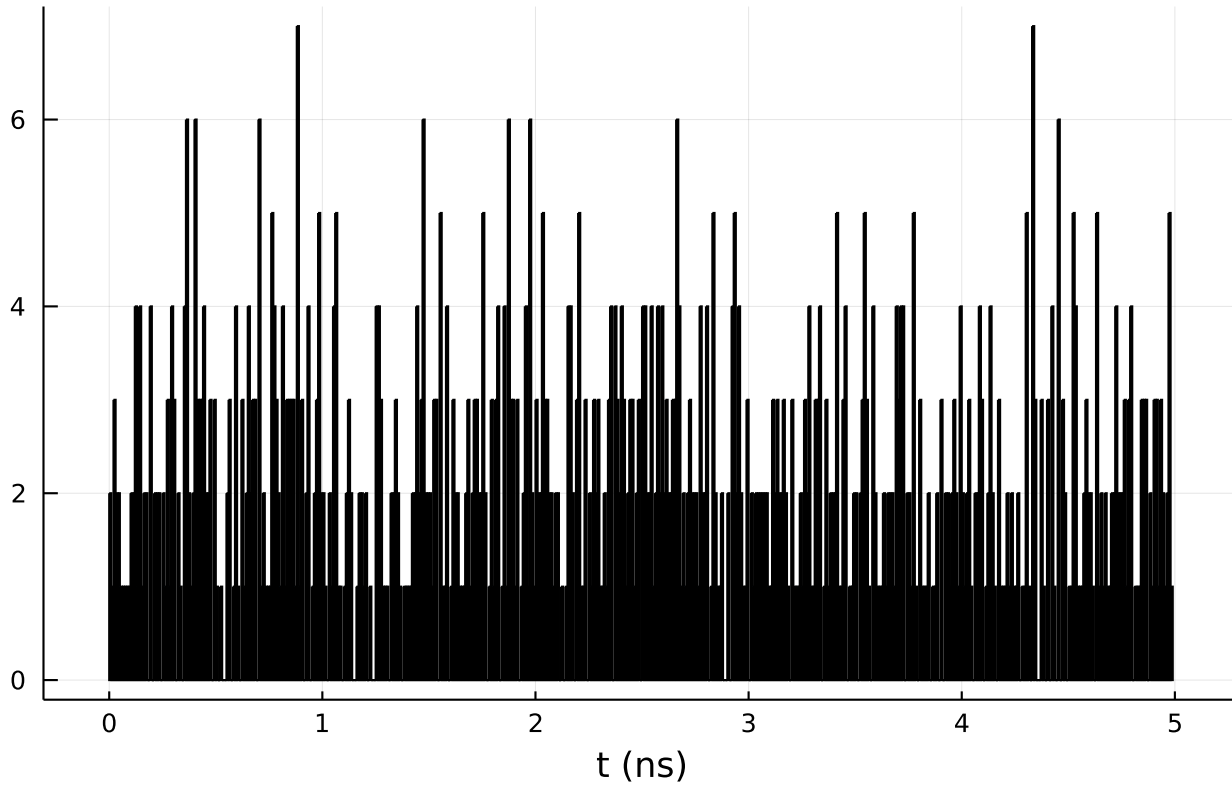


E-field/intensity: [Download...](#) classical_field_avg.csv

Correlations: [Download...](#) classical_correlations_avg.csv

Total counts:

Total counts displayed = 997



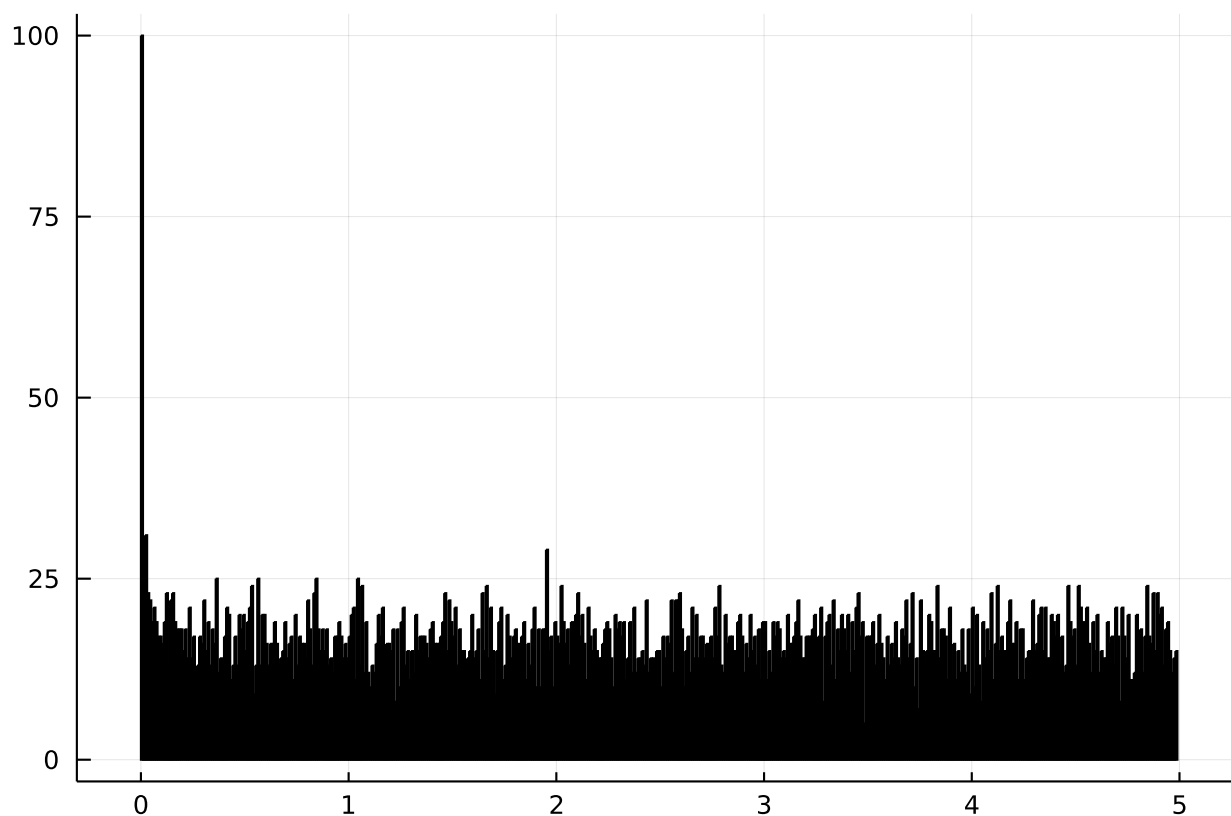
Bin edges: [Download...](#) photon-count_histogram_bin-edges.csv

Bin weights: [Download...](#) photon-count_histogram_bin-weights.csv

Arrival times: [Download...](#) photon-arrival-time_data.csv

Autocorrelation of photon counts

Note that the $\tau = 0$ bin matches the number of simulations. This only happens because we are using the autocorrelation for a single beam. This will not be true when I add a beam splitter.



Bin edges: [Download...](#) photon-correlation_histogram_bin-edges.csv

Bin weights: [Download...](#) photon-correlation_histogram_bin-weights.csv

Correlation times: [Download...](#) photon-correlation-tau_data.csv

Calculations

Classical calculations

```

• # these parameters are needed for all simulations
• begin
•     # Balmer-α lines specified here
•     ωM = 2*π*[456811.0, 456812.0]
•     # magnitude of each line
•     mag = convert(Vector{ComplexF64},ones(length(ωM)))
•     # calculate line differences
•     ΔM = ωM .- ωM[1]
•     # generate times in tres ps intervals up to 2*tmax
•     times = collect(0:tres*1e-3:2*tmax);
•     # limit the window to tmax to avoid correlation cutoff
•     window = convert(Integer,floor(length(times)/2));
•     # τ is just the times up to our window
•     τ = times[1:window];
• end;

• if makePlots
•     # our calculated g2τ
•     g2τCalc = 1
•     Emag2 = real.(mag .* conj(mag))
•     Emag4 = Emag2 .* Emag2
•     sumEmag2 = sum(Emag2)
•     sumEmag4 = sum(Emag4)
•     term2 = sumEmag4/(bigN*sumEmag2^2)
•     g2τCalc -= term2

•     term3 = sum(Emag2 .* exp.(-im*ΔM .* transpose(τ)),dims=1)/sumEmag2
•     term3 = real.(term3 .* conj(term3))

•     kbOverMhC2 = 9.178e-14;
•     σ = sqrt(kbOverMhC2*temp)*ωM[1]
•     stauAvg = transpose(bigN .+ bigN*(bigN-1)*exp.(-σ^2*τ .^2))
•     term3 = term3 .* stauAvg/bigN^2

•     g2τCalc = g2τCalc .+ term3
• end;

```

- Our calculated version

$$g^{(2)}(\tau) = 1 - \frac{1}{N} \frac{\sum_{m=1}^M |\mathcal{E}_m|^4}{\left(\sum_{m=1}^M |\mathcal{E}_m|^2\right)^2} + \left| \frac{\sum_{m=1}^M |\mathcal{E}_m|^2 e^{-i\Delta_m \tau}}{\sum_{m=1}^M |\mathcal{E}_m|^2} \right|^2 \frac{\langle S(\tau) \rangle_\omega}{N^2}$$

```

• # if ONLY the classical plots are desired, then this calculates a single instance
of the classical fields and correlations
• if makePlots && (classicalPlots && !(classicalAvgPlots || countsPlot || corrPlot))
•   # generate N doppler broadened frequencies
•   ω1Doppler = ωnDoppler(ωM[1],bigN,temp)
•   # generate N*M random phases
•   φ1nm = 2*π*rand(Float64,(length(ωM),bigN))
•   # construct field parameter object
•   testParams1 = eFieldParams(mag,ΔM,ω1Doppler,φ1nm)
•
•   # calculate electric field vs time
•   e1fieldt = map(t->electricField(t,testParams1),times)
•
•   # calculate g1τ
•   g1τ1Norm = correlate(e1fieldt,conj.(e1fieldt),0>window);
•   g1τ1 = abs.(map(i->correlate(e1fieldt,conj.
(e1fieldt),i>window),collect(0>window-1))/g1τ1Norm)
•
•   # calculate the intensity vs time
•   intensity1t = real.( e1fieldt .* conj(e1fieldt))
•
•   # calculate g2τ
•   g2τ1Norm = mean(intensity1t)^2
•   g2τ1 = map(i->autocorrelate(intensity1t,i>window),collect(0>window-1))/g2τ1Norm
• end;

```



```

• # the classical average plots and photon-based plots require multiple instances of
the classical field calculations
• if makePlots && (classicalAvgPlots || countsPlot || corrPlot)
•   # calculate number of trials from the total desired photon counts and the
average photon count per trial
•   nTrials = convert(Integer,ceil(ntot/nbar))
•   # make an array of the average photon counts per trial for array broadcasting
•   nPerTrial = bigN*ones(Integer,nTrials)
•   # calculate nTrials instances of doppler broadened frequencies
•   ωDoppler = ωDoppler.(ωM[1],nPerTrial,temp);
•   # calculate nTrials instances of random phases
•   φnm = map(n->2*π*rand(Float64,(length(ωM),n)),nPerTrial);
•   # generate nTrials instances of field parameters
•   testParams = map((x,y)->eFieldParams(mag,ΔM,x,y),ωDoppler,φnm);
•
•   # calculate nTrials instances of the time dependent electric field
•   efieldt = map(x->map(t->electricField(t,x),times),testParams);
•
•   # calculate nTrials instances of g1τ
•   g1τNorm = map(eft->correlate(eft,conj.(eft),0>window),efieldt);
•   g1τ = map((eft,normG1τ)->abs.(map(i->correlate(eft,conj.
(eft),i>window),collect(0>window-1))/normG1τ),efieldt,g1τNorm);
•
•   # calculate nTrials instances of the time dependent intensity
•   intensityt = map(eft->real.( eft .* conj(eft)),efieldt);
•
•   # calculate nTrials instances of g2τ
•   g2τNorm = map(intens->mean(intens)^2,intensityt)
•   g2τ = map((intens,normG2τ)->map(i-
>autocorrelate(intens,i>window),collect(0>window-1))/normG2τ,intensityt,g2τNorm)
•
•   # pick out one instance of everything for plotting
•   e1fieldtM = efieldt[1]
•   g1τ1M = g1τ[1]
•   intensity1tM = intensityt[1]
•   g2τ1M = g2τ[1]
• end;

```

```

• if makePlots && classicalAvgPlots
•   efieldtAvg = vectorAvg(efieldt)
•   g1τAvg = vectorAvg(g1τ)
•   intensitytAvg = vectorAvg(intensityt)
•   g2τAvg = vectorAvg(g2τ)
• end;

```

Calculate photon counts

Photon counts are calculated by treating the intensity in each time bin as the average photon count rate, then sampling from a poisson distribution with that average count rate.

```

• if makePlots && (countsPlot || corrPlot)
•     yCounts = map(intens->poissonCount.(yIntensity(intens,nbar*2)),intensityt)
•     ycountTimes = map(yCt->countTimes(times,yCt),yCounts)
•     flatyCountTimes = vcat(ycountTimes...)
• end;

```

Calculate autocorrelation of photon counts

Note that I only look at whether two bins both have counts or not when calculating the autocorrelation. I **do not** look at *how many* counts there are in each bin.

```

• if makePlots && corrPlot
•     correlationTimes = map(yCt->singleDeltaTimes( $\tau$ ,yCt),yCounts)
•     flatCorrelationTimes = vcat(correlationTimes...)
• end;

```

Functions

Main.workspace2558.singleDeltaTimes

```

• """
•     function singleDeltaTimes( $\tau$ ::Vector,yCounts::Vector)
•
•     Returns an array of  $\tau$  values for which the yCounts autocorrelation is non-zero.
•     """
•     function singleDeltaTimes( $\tau$ ::Vector,yCounts::Vector)
•         return  $\tau$ [map(i->autocorrelate(yCounts,i,length( $\tau$ )) > 0 ? true :
•             false,collect(0:length( $\tau$ )-1)) ]
•     end
•

```

Main.workspace2568.countTimes

```

    """
    function countDeltaTimes( $\tau$ ::Vector,  $\gamma$ Counts::Vector)
    Returns an array of  $\tau$  values for which the  $\gamma$ Counts autocorrelation is non-zero.
    """
    function countTimes(times::Vector,  $\gamma$ Counts::Vector)
        out = Vector{Real}(undef, 0)
        for (i, counts) in enumerate( $\gamma$ Counts)
            if counts != 0
                countTimes = times[i]*ones(counts)
                out = vcat(out, countTimes)
            end
        end
        return out
    end
end

```

Main.workspace3.eFieldParams

```

    """
    eFieldParams64(mag::Array{S},  $\Delta$ M::Array{T},  $\omega$ N::Array{T},  $\phi$ ::Array{S} ) where
    {T<:Real, S<:Complex}
    Static parameters for the electric field
    """
    struct eFieldParams
        mag::Vector
         $\Delta$ M::Vector
         $\omega$ N::Vector
         $\phi$ ::Matrix

        function eFieldParams(mag::Vector,  $\Delta$ M::Vector,  $\omega$ N::Vector,  $\phi$ ::Matrix )
            @assert length(mag) == length( $\Delta$ M) "Number of magnitudes must match number
            of emission lines"
            @assert (size( $\phi$ )[1] == length( $\Delta$ M) && size( $\phi$ )[2] == length( $\omega$ N)) "Must have a
            unique phase for each n and m"
            new(
                convert(Vector{Complex}, mag),
                convert(Vector{Real},  $\Delta$ M),
                convert(Vector{Real},  $\omega$ N),
                convert(Matrix{Real},  $\phi$ )
            )
        end
    end
end

```

Main.workspace3.electricField

```

    """
    function electricField(t::Real,params::eFieldParams)
    Returns the electric field value at time t
    """
    function electricField(t::Real,params::eFieldParams)
        # generate frequencies
        ωNM = transpose(params.ωN) .+ params.ΔM
        # add the phase
        exponentnm = -im*(t*ωNM+params.φ)
        # put them in the exponent
        enm = exp.(exponentnm)
        # multiply by the field magnitude
        fieldnm = params.mag .* enm
        # add it all together
        return sum(ivec(fieldnm))
    end

```

Main.workspace3.ωNDoppler

```

    """
    function ωNDoppler(ω0::Real,N::Integer,temp::Real,seed::Integer = -1)
    Generates N doppler shifted frequencies around frequency ω0 for a source at
    temperature temp. Seed optional for reproducible results.
    """
    function ωNDoppler(ω0::Real,N::Integer,temp::Real,seed::Integer = -1)
        rng = MersenneTwister()
        if seed != -1
            rng = MersenneTwister(seed)
        end
        kbOverMhC2 = 9.178e-14;
        σ = sqrt(kbOverMhC2*temp)*ω0
        d = Normal(ω0,σ)
        return rand(rng,d,N)
    end

```

Main.workspace3.correlate

```

"""
    function correlate(u::Vector{T},v::Vector{T},offset::Integer>window::Integer =
-1) where {T<:Number}
    Calculates correlation between vectors u and v with given offset. Specify averaging
    window to limit range of correlation. If the window extends beyond the end of one
    vector, it treats out-of-bounds indices as zero.
"""
function correlate(u::Vector{T},v::Vector{T},offset::Integer>window::Integer = -1)
where {T<:Number}
    @assert offset <= length(u) "Offset out of bounds"
    @assert window <= length(u) && window <= length(v) "Window must be smaller than
input vector lengths"
    if window == -1
        window = length(u)
    end
    v1 = view(u,1:window)
    v2 = view(v,1+offset:min(window+offset,length(v)))
    if window+offset > length(v)
        v2 = vcat(v2,zeros(window+offset-length(v)))
    end
    return dot(v1,v2)/window
end

```

Main.workspace3.autocorrelate

```

"""
    function autocorrelate(u::Vector{T},offset::Integer, window::Integer = -1)
where {T<:Number}
    Calculates correlation of vector u with itself.
"""
function autocorrelate(u::Vector{T},offset::Integer, window::Integer = -1) where
{T<:Number}
    correlate(u,u,offset>window)
end

```

Main.workspace3.γIntensity

```

"""
    function γIntensity(intensity::Vector,nbar::Real)
    Calculates the photon count rate in each bin of an intensity histogram
"""
function γIntensity(intensity::Vector,nbar::Real)
    nintensity = intensity/sum(intensity)
    return nbar*nintensity
end

```

Main.workspace3.poissonCount

```

    """
    function poissonCount(nbar::Real)
    Returns Poisson distributed counts for average count rate nbar
    """
    function poissonCount(nbar::Real)
        p = exp(-nbar)
        s = p
        r = rand()
        count = 0
        while r > s
            count += 1
            p *= nbar/count
            s += p
        end
        return count
    end
end

```

Main.workspace3.beCount

```

    """
    function beCount(nbar::Real)
    Returns Bose-Einstein distributed counts for average count rate nbar
    """
    function beCount(nbar::Real)
        p = 1/(nbar+1)
        fnbar = p*nbar
        f = p*nbar
        s = p
        r = rand()
        count = 0
        while r>s
            count += 1
            p *= f
            s += p
        end
        return count
    end
end

```

vectorAvg (generic function with 1 method)

```

    function vectorAvg(someVector::Vector)
        return +(someVector...)/length(someVector)
    end

```

Load Prerequisites

```
• import Pkg

• Pkg.add("Distributions")

• Pkg.add("DataFrames")

• Pkg.add("CSV")

• Pkg.add("Plots")

• Pkg.add("IterTools")

• Pkg.add("PlutoUI")

• Pkg.add("StatsBase")

• using Random, Distributions, StatsBase

• using LinearAlgebra

• using Plots

• using IterTools

• using PlutoUI

• using DataFrames

• using CSV
```

