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Liquid xenon scintillation measurements and pulse shape discrimination in the LUX dark matter detector

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Weakly Interacting Massive Particles (WIMPs) are a leading candidate for dark matter and are expected to produce nuclear recoil (NR) events within liquid xenon time-projection chambers. We present a measurement of liquid xenon scintillation characteristics in the LUX dark matter detector and develop a pulse shaped based discrimination parameter to be used for particle identification. To accurately measure the scintillation characteristics, we develop a template-fitting method to reconstruct the detection time of photons. Analyzing calibration data collected during the 2013-16 LUX WIMP search, we measure a singlet-to-triplet scintillation ratio for electron recoils (ER) that is consistent with existing literature, and we make a first-ever measurement of the NR singlet-totriplet ratio at recoil energies below 74 keV. A prompt fraction discrimination parameter exploits the difference of the photon time spectra for NR and ER events and is optimized to have the least number of ER events that occur in the 50% NR acceptance region. When this discriminator is used in conjunction with charge-to-light discrimination on the calibration data, the signal-to-noise ratio in the NR dark matter acceptance region increases by up to a factor of two.

I. INTRODUCTION

Liquid xenon time projection chamber (TPC) experiments are leaders in sensitivity to detect the interactions of Weakly Interacting Massive Particles (WIMPs), a class of as-yet-unobserved particles that have been proposed as a solution to the dark matter problem [\[1](#page-15-0)[–3\]](#page-15-1). In such experiments, the WIMP is predicted to scatter elastically from a xenon nucleus, resulting in a nuclear recoil (NR). The primary backgrounds are electron recoils (ER) from gamma and beta radiation released by residual radioactivity in the detector materials, with a small contribution from neutrino-electron scattering. Interactions in liquid xenon produce scintillation photons and ionization electrons which can be measured to reconstruct information about the interaction. TPC experiments measure both the ionization and scintillation signals and use this information to reconstruct the energy deposition, particle type, and the position of the interaction.

Background rejection is paramount to the success of liquid xenon dark matter searches. Material screening and shielding are the primary methods to mitigate backgrounds; detectors are constructed from highly radiopure materials and are operated in well-shielded underground environments to reduce backgrounds from cosmic rays and environmental sources. Position reconstruction allows fiducialization and the rejection of multiplescattering events. The former eliminates ER backgrounds from detector materials stopping close the edges of the sensitive volume, while the latter removes event topologies inconsistent with WIMP scattering. Background events which remain in the active volume can be rejected through particle-type discrimination between ER and NR [\[1\]](#page-15-0). In liquid xenon TPC experiments, this last step is typically done using the ratio of ionization charge to scintillation light in the event, which is higher for ER events than NR events. The present work explores enhancing the ER background rejection using pulse shape discrimination (PSD) applied to the scintillation signal alone.

Scintillation light is produced by the self-trapping of excited xenon atoms (Xe[∗]), created when a particle deposits energy in the liquid. Direct excitation and recombination of electron–ion pairs create excited atoms, which combine with a neutral ground-state Xe atom to form the molecular dimer Xe_2^* . The dimer decays to the monatomic ground state via emission of a VUV photon $(\lambda = 175 \text{ nm})$ [\[4,](#page-15-2) [5\]](#page-15-3). These two processes are shown

schematically in Eq. [1.1](#page-2-2) (direct excitation) and Eq. [1.2](#page-2-3) (recombination of electron–ion pair).

$$
Xe^* + Xe \rightarrow Xe_2^* \n\rightarrow Xe + Xe + \gamma,
$$
\n(1.1)

$$
Xe^{+} + Xe \rightarrow Xe_{2}^{+}
$$

\n
$$
Xe_{2}^{+} + e^{-} \rightarrow Xe^{**} + Xe
$$

\n
$$
Xe^{*} + Xe \rightarrow Xe_{2}^{*}
$$

\n
$$
\rightarrow Xe + Xe + \gamma.
$$

\n(1.2)

The decay of the dimer is observed to have both a fast and a slow component, which are interpreted as de-excitation of the singlet ${}^{1}\Sigma_{u}^{+}$ and the triplet ${}^{3}\Sigma_{u}^{+}$ states, respectively [\[6,](#page-15-4) [7\]](#page-15-5). There are conflicting measurements of the lifetimes of these states in the literature; measurements of the singlet time constant τ_1 range from 2 to 4 ns, while measurements of the triplet time constant τ_3 range from 21 to 28 ns [\[7,](#page-15-5) [8\]](#page-15-6). Both components have been observed for electron recoils, alpha recoils, and recoiling fission fragments.

For electron recoils, some experiments operating without an applied electric field have observed a time profile that is best fit with a single exponential with $\tau = 30$ – 45 ns [\[7,](#page-15-5) [9,](#page-15-7) [10\]](#page-15-8). This is attributed to an additional time delay due to electron–ion recombination. This interpretation is supported by measurements that show that the scintillation time structure reduces to the characteristic singlet/triplet shape under an applied electric field (which suppresses recombination) [\[6\]](#page-15-4). In addition, recent measurements, without an applied field, show an energy-dependence of the long component, correlated with the energy-dependence of recombination [\[11\]](#page-15-9). No field-dependence is observed for alpha particle or fission fragment recoils, suggesting that recombination-related timing effects are only significant at low ionization densities. At the energies (0-50 keV) and electric fields (100– 1000 V/cm) relevant for modern liquid xenon TPC experiments, there are no direct measurements of the effects of recombination on ER scintillation timing. However, extrapolating to this regime using the empirical model given in Ref. [\[12\]](#page-16-0) suggests that recombination does not play a significant role in scintillation emission timing in these experiments, and that pulse shapes can be welldescribed purely in terms of the singlet and triplet emission.

The ratio of singlet emission to triplet emission varies with particle type, opening up the possibility for ER/NR discrimination using PSD. Multiple groups have studied liquid xenon PSD in small R&D detectors [\[10,](#page-15-8) [13,](#page-16-1) [14\]](#page-16-2), and it was successfully used to reduce backgrounds in early liquid-xenon-based dark matter searches [\[15–](#page-16-3)[18\]](#page-16-4).

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However, these studies are restricted to small detectors or detectors with spherical photosensor coverage of the xenon volume. Current and future TPC experiments have meter-scale dimensions and make extensive use of reflectors to maximize light collection [\[19,](#page-16-5) [20\]](#page-16-6). In such detectors, scintillation pulse shapes are subject to significant distortion from scattering, reflection, and absorption of photons by detector materials. In addition, previous studies have not attempted to reconstruct the singlet/triplet ratio for both ER and NR pulses at the low energies relevant to dark matter searches. Attempts to simulate scintillation pulses must therefore rely on measurements at higher energies, which may not accurately reflect xenon microphysics in the region of interest.

In this work, we present a measurement of scintillation characteristics and PSD in the LUX detector, a meterscale liquid xenon TPC [\[21\]](#page-16-7). We study both ER and NR calibration data taken throughout the LUX WIMPsearch campaign. First, a template-based photon reconstruction algorithm is used to deconvolve the response of the electronics in order to reconstruct the time when a photon strikes a photomultiplier tube. The spectra of photon detection times are added across many pulses to construct average pulse shapes for both ER and NR events. Second, we develop an analytical model to decouple detector effects from xenon scintillation emission. This model is fit to data to extract physical parameters that can inform simulation packages such as NEST [\[22\]](#page-16-8). Finally, we construct a pulse shape discriminant using the prompt-fraction technique, and compute the power of PSD background rejection in LUX. Using the bestfit parameters from the analytical model, we construct a simulation that accurately reproduces PSD measurements from data. The discrimination power improves with recoil energy, and we demonstrate how PSD can be used in conjunction with the charge-to-light ratio to further improve background rejection. These features make it attractive for exotic dark matter searches in which low-energy recoils are suppressed, such as searches for momentum-dependent and inelastic dark matter scattering from nuclei [\[23,](#page-16-9) [24\]](#page-16-10). These measurements allow estimation the of the PSD capabilities of the current and next generation of liquid xenon dark matter experiments, and can be applied to future dark matter searches using the LUX dataset.

II. THE LUX EXPERIMENT

The LUX detector is a dual-phase xenon TPC, designed to detect WIMP scattering with xenon nuclei. It was operated from April 2013 through June 2016 in the Davis Cavern at the Sanford Underground Research Facility (SURF) in Lead, South Dakota [\[25\]](#page-16-11). Dark matter search data were acquired in two exposure periods, denoted WS2013 and WS2014-16 [\[1\]](#page-15-0). To meet the stringent low-background requirements required for the dark matter search, the detector is located deep underground

(4,300 meters water equivalent), is surrounded by a 7.6 m tall by 6.1 m diameter water shield, and is constructed from materials that have been carefully screened for radiopurity. The sensitive volume is approximately 48 cm in height and 24 cm in radius, and contains ∼250 kg of liquid xenon. Each end of the TPC is instrumented with an array of 61 Hamamatsu R8778 photomultipler tubes (PMTs) to detect light signals generated in the TPC. Twelve PTFE panels, >95% reflective at 175 nm [\[26\]](#page-16-12), line the walls to increase the light collection efficiency. The scintillation signal, denoted S1, is detected directly by the PMTs. Ionization electrons are drifted under an applied electric field and extracted into a gas region at the top of the detector, producing an electroluminescence signal, denoted S2. The (x, y) position of the events is reconstructed using the pattern of S2 light on the top PMT array [\[27\]](#page-16-13), while the depth is reconstructed from the time delay between the S1 and S2 signals. The energy deposition of the event is reconstructed from the magnitude of the two signals.

The PMT signals are routed to an external electronics breakout box for processing, before digitization. Signals are amplified in two stages at the pre- and postamplifiers, which provide a total effective gain of 7.5. The signals are shaped by a 30 MHz low-pass filter. The resulting single photoelectron (SPE) pulses have a full width at half maximum (FWHM) of >20 ns [\[28\]](#page-16-14). Signals are digitized using a 100 MHz Struck SIS3301 8-channel fast analog to digital converter (ADC). The average digitized area of an SPE is ∼100 mV·ns [\[21\]](#page-16-7). Recently, it has been shown that VUV photons have a ∼20% probability of generating two photoelectrons at the photocathode of the R8778 PMT [\[29\]](#page-16-15). Pulse areas are therefore measured in units of detected photons (phd) rather than photoelectrons. In addition, a "spike count" has been employed in the LUX dark matter analyses to improve resolution for very low-energy S1 pulses [\[1,](#page-15-0) [30\]](#page-16-16). This method is not used in the present work, as we are focused on higher energy events.

Calibration campaigns were conducted throughout the exposure period to monitor detector stability and response. Detector stability, electron lifetime, and signal corrections were measured using a 83mKr source, dissolved in xenon [\[31\]](#page-16-17). The corrected signals for the S1 and S2 pulses are denoted S1 and S2. These calibrations occurred weekly throughout both exposure periods. The low-energy NR and ER responses were periodically calibrated in–situ using fast neutrons from deuteriumdeuterium fusion (DD) [\[32\]](#page-16-18) and electrons from the beta decays of tritium [\[33\]](#page-16-19). Neutrons from the DD generator traveled through an air-filled collimating pipe suspended in the water tank, and were approximately normally incident on the detector at a level ∼7.5 cm below the LXe surface in the TPC. The tritium source was deployed in the form of tritiated methane CH3T and mixed into the detector through the xenon circulation system. Both DD and CH₃T calibrations were performed at the end of each WIMP search run, as well as three times during WS2014-16. To calibrate the depth-dependent response to NR events, DD calibrations were conducted at different heights at the start and end of WS2014-16. Additional calibration campaigns were carried out at the end of WS2014-16, including an injection of ¹⁴C into the xenon circulation system. The higher energy beta spectrum provided by ${}^{14}C$ (endpoint at 156 keV) provides a source of ER events beyond the 18 keV endpoint of tritium.

We use all of the DD, CH_3T , and ¹⁴C calibration data in the analysis presented here. Due to limited statistics in the lower portion of the detector, we developed our analysis and fit our analytical model using the data in the top drift bin from the WS2014-16 WIMP analysis $(t_{drift} = 40{\text -}105 \,\mu s)$ and demonstrate consistency between data and simulations in the lower drift bins.

III. PHOTON TIMING

In past studies, the time structure of detected scintillation photons was typically obtained by measuring the shape of pulses summed over all channels in a detector. However, the ∼20 ns shaping time constant and 10 ns sampling period of the LUX DAQ are similar to the timescale to the de-excitation process of liquid xenon, and may therefore mask underlying scintillation characteristics. For this analysis, we developed a photon timing algorithm and a channel-to-channel time calibration technique that accurately reconstructs a photon's detection time by deconvolving it from the electronic pulse.

A. Photon timing algorithm

Precise timing is achieved with an analysis technique that separates pulses into individual detected photons, similar to the approaches in Refs. [\[11,](#page-15-9) [34\]](#page-16-20). After baseline subtraction and normalization by PMT gains, the waveforms in individual PMT channels were analyzed in three steps: 1) template model fitting, 2) template model selection, and 3) re-weighting of the reconstructed photons.

In the template fitting stage, the waveform in a single channel is fit with five separate n -photon models, composed of the sum of up to five single-photon template functions (the restriction to $n \leq 5$ is expected to be more than 99% efficient for scintillation pulses up to 300 detected photons). The single-photon template is an empirical model, constructed from an average of 1,000 waveforms with areas between 0.5–1.5 phd. The fit is performed using the Migrad routine built into the TMinuit class in ROOT [\[35\]](#page-16-21), with the amplitude and the arrival time of each template as free parameters. Initial values for times and amplitudes are given by the time and height of the peaks in the waveform, defined as maxima above a threshold of 0.1 phd/sample ($\sim 5\sigma$ above baseline fluctuations). When there are fewer than five peaks, the photon fit is repeated with all possible permutations, allowing multiple detected photons piling up to form a single spike. The resulting fits must meet two criteria: none of the reconstructed photons may have an area less than 0.15 phd, and the time separation between all pairs of photons must be greater than one sample. The first criterion removes fits in which we reconstruct fluctuations in baseline noise. Roughly 2% of real photons fall below this threshold. The second criterion removes fits with multiple photons reconstructed within a single sample, where our algorithm is unable to accurately separate photons (a correction is applied later to account for unresolved pileup). If a particular fit fails one or both of these cuts. The best-fit times, areas, and likelihood values for each of the remaining n-photon models are passed to the next stage of the algorithm for comparison and selection.

The model comparison stage uses Bayes' Theorem to assign a likelihood score to each n-photon model, and the model with the highest likelihood score is selected. The likelihood score is the product of the maximum likelihood from the fit and a prior probability calculated using the measured area. For a given waveform (denoted D) and n-photon model (denoted \mathcal{M}_n), Bayes' Theorem can be used to calculate the probability of \mathcal{M}_n given D, with,

$$
P(\mathcal{M}_n|D) = \frac{P(D|\mathcal{M}_n)P(\mathcal{M}_n)}{P(D)}.
$$
 (3.1)

Here, $P(D|M_n)$ is the maximum likelihood given by the fit, $P(D)$ is a flat normalization constant which we ignore, and $P(\mathcal{M}_n)$ is the prior. The prior $P(\mathcal{M}_n)$, is the probability of measuring the observed area if there were actually n detected photons in the channel. This is calculated using a single-photon area response PDF, averaged over all PMT channels, which incorporates the $\sim 20\%$ probability of xenon scintillation light producing two photoelectrons in the R8778 PMTs [\[29\]](#page-16-15). The $P(\mathcal{M}_n)$ prior prevents overfitting by applying a penalty to models composed of many reconstructed photons with improbably small areas. The model with the largest overall likelihood score $P(\mathcal{M}_n|D)$ is selected and we return the best-fit arrival times and amplitudes.

To correct for unresolved pile-up, we assign a weight to each reconstructed photon that is equal to the area of the fitted template. Pile-up occurs when the fitting algorithm fails to split a single peak, usually when two photons arrive in a single channel within 1 sample. The total number of photons counted in the pulse is given by,

Corrected photon count =
$$
C \sum_{i=1}^{N} w_i
$$
, (3.2)

where w_i is the weight of the *i*-th photon, N is the uncorrected number of photons returned by the fits in all channels, and C is an overall correction factor. The latter accounts for inefficiencies that may arise due to the fit threshold or inexact area matching of the template function with true pulses. To find C , we fit a linear model of the form $y = mx$ to the DD neutron data, where x is the pulse area (in phd) and $y = \sum w_i$. Then $C = 1/m$. We

FIG. 1: Comparison of corrected photon count vs. the raw S1 pulse area in detected photons (phd) for both tritium (blue) and DD neutron (red) calibration data. The corrected photon count is computed using Eq. [3.2](#page-4-0) with $C = 1.04$. The black dashed line has a slope $m = 1$, and is shown for comparison.

find that $C = 1.04 \pm 0.01$ reproduces the total number of photons obtained from the pulse area. A comparison between our corrected photon count and the pulse area using these values is shown in Fig. [1.](#page-5-0) The pulse area and the corrected photon count agree throughout the 0– 200 phd pulse area range used in this work.

The times of the photons returned by the fits correspond to the photon detection times, deconvolved from the shaping of the detector electronics. The performance of our algorithm is demonstrated in Figs. [2](#page-5-1) and [3.](#page-6-0) Figure [2](#page-5-1) shows an example best-fit model with a simulated pulse. Although there are only two peaks in the pulse, the model selection algorithm correctly prefers the 3-photon model and reconstructs the times to within 0.1 samples (1 ns) for each photon. Fig. [3](#page-6-0) shows the algorithm applied to a real S1 pulse from the tritium calibration data. We estimate the 1σ uncertainty of the photon detection time from the fit to be 1.6 ns, calculated from the average uncertainty returned by the Migrad fitter when our algorithm is applied to the CH_3 T calibration data.

B. Channel-to-Channel Time Calibration

There are several factors that affect the relative time between PMT channels. The R8778 PMTs are specified to have an electron transit time of 41 ± 1.7 (σ) ns at 1500 V; this transit time varies inversely with the square root of the bias voltage [\[36\]](#page-16-22). Gain-matching of the PMTs requires operational voltages to vary between 1000 and 1500 V, which causes the electron transit time to vary be-

FIG. 2: Example simulated waveform (black). Three photons arriving at $t = -0.75, 1.60,$ and 6.50 samples $(1 \text{ sample} = 10 \text{ ns})$ are used to generate the simulated signal. The photon timing algorithm described in Section [III](#page-4-1) reconstructs 3 photons arriving at $t =$ −0.76, 1.51, and 6.45 samples (grey).

tween 41 and 50 ns. Differences in cable lengths can cause further differences in signal arrival times. The shaping filters on the pre- and post-amplifiers further degrade timing accuracy and can add relative delays between channels. Finally, a 100 MHz clock pulse is propagated to each digitizer that can cause synchronization delays between digitizers [\[21\]](#page-16-7). All of these relative offsets must be measured and corrected for in this analysis so that coincident photons are correctly aligned in time.

We measure the combined effect of these time offsets using LEDs mounted on the top and bottom PMT arrays. The LED system includes twelve 440 nm diodes, capped with PTFE diffusers, that are used for gain and afterpulsing calibrations of the PMTs. To measure relative time offsets, pulses with a FWHM of 20 ns, a rise/falltime of 5 ns, and a peak amplitude in the range 3.36– 3.80 V are propagated to individual LEDs within the chamber. The resulting single photoelectron pulses in each channel are fit with the single-photon template to determine their arrival time relative to the LED strobe. The direct-path travel time for a photon from an LED to a PMT is subtracted to remove photon path length differences from this calibration. If the PMT is located on the same array as the LED, there is no direct optical path from the LED to the PMT; we therefore assume that the shortest path is via reflection off the liquid-gas interface. For each channel, a distribution of the path-corrected photon detection time, relative to the LED trigger, is obtained. A typical distribution is shown in Fig. [4.](#page-6-1)

A common reference time needs to be selected for each channel to serve as the correction to be applied to that channel. Tails longer than the strobe time of the LED are

(a) Measured Scintillation Pulse

(b) Reconstructed Photon Times

FIG. 3: (a) A scintillation pulse from a CH_3T calibration event, separated by PMT channel. (b) The reconstructed peak times for the photons in the pulse is shown.

observed in these time distributions and are attributed to photons scattering within the detector volume. To avoid these effects from biasing our measurement, 10% of the peak amplitude on the rising edge of each distribution is used to define the correction time for each channel. These corrections vary by up to 20 ns (2 samples) from channel to channel. Measurements are repeated with two LEDs in the top array and two LEDs in the bottom array to test for systematic effects from PMT coverage. The corrections between different LED measurements agree to within 2 ns, which represents the resolution of the measurement; we take this as our 1σ uncertainty. The corrections are subtracted from the reconstructed photon

FIG. 4: An example of a photon detection time distribution for a PMT located on the bottom array, facing the strobed LED on the top array. The photon detection times shown are relative to the LED trigger and are histogramed in 2 ns bins. The points between the data are interpolated with a linear spline, which is shown with the dotted red trace. The dashed blue line shows 10% of the peak height and is compared to the spline fit to obtain the offset for that channel, shown by the solid green trace.

times in the analysis presented in Section [V.](#page-8-0)

IV. LIQUID XENON SCINTILLATION IN THE LUX DETECTOR

A. Analytical model of photon detection times

The analytical model of photon detection time in LUX is built from three components: scintillation emission, optical transport, and a model of instrument response.

The scintillation emission distribution is assumed to be of the form

$$
P_s(t) = C_1 e^{-t/\tau_1} + C_3 e^{-t/\tau_3}, \qquad (4.1)
$$

where τ_1 and τ_3 are the time constants governing the decay of the singlet and triplet states. In this parameterization, the ratio of singlet photons to triplet photons, referred to as the intensity ratio or the singlet/triplet ratio in the literature, is given by $(C_1 \tau_1) / (C_3 \tau_3)$. Additional timing effects in electron recoils due to electron– ion recombination are neglected in our model as they are expected to be suppressed by the applied electric field and the high linear energy transfer (LET) at low energies. For the calibration data used to constrain the model (\sim 300 V/cm and \sim 10 keV), an empirical formula in Ref. [\[12\]](#page-16-0) predicts a recombination time scale of τ_{rec} < 0.7 ns. As this is significantly smaller than the

TABLE I: Optical transport parameters at different locations inside the LUX detector (see Eq. [4.2\)](#page-7-0). The first four rows correspond to the drift bins used in the WS2014-16 dark matter analysis [\[1\]](#page-15-0), while the D-D beam location is used to constrain our pulse shape model in Sec. [V B.](#page-9-0)

| Position | \overline{A} | B_a | τ_a (ns) | τ_b (ns) |
|---------------|----------------|-------|---------------|---------------|
| Top | 0.0544 | 1.059 | 11.2 | 2.80 |
| Top center | 0.0489 | 1.017 | 11.2 | 5.21 |
| Bottom center | 0.0586 | 0.906 | 11.2 | 1.56 |
| Bottom | 0.120 | 0.798 | 11.1 | 1.68 |
| D-D beam | 0.0574 | 1.062 | 11 1 | 2.70 |

FIG. 5: Simulated time distribution of photons detected in the LUX detector, calculated using the ray-tracing capabilities in the LUXSim Geant4 simulation package. The photon sources are 10 cm thick, 20 cm diameter cylinders, centered at different depths below the liquid surface: 42 cm (grey), 32 cm (green), 22 cm (orange), and 12 cm (blue). These slices in depth correspond to the drift time bins used in the WS2014-16 dark matter search [\[1\]](#page-15-0). Solid lines show the second term in Eq. [4.2,](#page-7-0) fitted to these simulations. The excess in the first bin from direct-path photon arrivals is parameterized by a constant A and is not reflected in the curves shown. Values of the parameters for each position are given in Table [I](#page-7-1)

other timescales in this analysis, we neglect a full treatment and simply use two different triplet time constant for ER and NR, $\tau_{3_{er}}$ and $\tau_{3_{nr}}$. Any recombination effects will be absorbed by $\tau_{3_{er}}$ and will result in $\tau_{3_{er}}$ slightly larger than $\tau_{3_{nr}}$.

The optical transport distribution is constructed using photon-tracking simulations which take into account physical and geometrical effects on xenon scintillation light inside the LUX detector [\[37\]](#page-16-23). Photon transport times depend on several physical properties of the detector internals: reflection and absorption at internal surfaces, reflection and transmission at the liquid/gas interface, and absorption and scattering in the liquid. The values for these parameters are constrained in-situ using $83m$ Kr calibration data [\[38\]](#page-16-24). Photon transport is simulated using the LUXSim package [\[39\]](#page-16-25), a Monte Carlo code built on Geant4 [\[40,](#page-16-26) [41\]](#page-16-27). Figure [5](#page-7-2) shows the optical transport distributions in each of four height bins used in the LUX WIMP search analysis with the WS2014-16 exposure [\[1\]](#page-15-0). The differences between the simulations reflect the depth-dependent probability, due to the combination of geometric efficiency and reflection at the liquidgas interface, for photons to travel directly to PMTs. To include optical transport in the analytical model, we introduce the empirical distribution

$$
P_o(t) = A \,\delta(t) + (1 - A) \left[\frac{B_a}{\tau_a} \, e^{-t/\tau_a} + \frac{B_b}{\tau_b} \, e^{-t/\tau_b} \right],\tag{4.2}
$$

where A, B_a , τ_a , and τ_b are fitted to the simulated distributions. The first term is a Dirac delta function which parameterizes the light which travels directly to a PMT, while the second term parameterizes the time distribution from light that reflects and scatters from the detector internals. Normalization requires $B_b = 1 - B_a$. The uncertainties in the optical parameters, given in Ref. [\[38\]](#page-16-24), are a source of systematic error in our analysis. We discuss this further in Section [V B.](#page-9-0)

Finally, we treat instrumental effects as normally distributed variables, parameterized by an overall width σ . There are two leading sources of random timing fluctuations in our data: the transit time spread of the R8778 PMTs $\sigma_{tts} = 1.9$ ns (at the operating bias of ∼1300 V) [\[42\]](#page-16-28), and the uncertainty in the reconstructed detection time from the photon timing algorithm, $\sigma_{fit} = 1.6$ ns. There is also the 2 ns uncertainty in the channel-tochannel time corrections, σ_{tc} . While this is a fixed time offset for each channel rather than a random pulse-bypulse fluctuation, the result is a net smearing of the pulse shape when averaging pulses together across all of the channels. Using simulations, the net effect of the time correction uncertainties on average pulse shapes was determined to be equivalent to normally-distributed random fluctuations. Therefore, the overall width added to photon time spectra by the effects in the electronics and the data reduction pipeline can be described by adding these three effects in quadrature:

$$
\sigma = \sqrt{\sigma_{tts}^2 + \sigma_{tc}^2 + \sigma_{fit}^2} \,. \tag{4.3}
$$

The total analytical model for photon detection time is given by the convolution of Eq. [4.1,](#page-6-2) Eq. [4.2,](#page-7-0) and a Gaussian distribution of width σ , given by Eq. [4.3:](#page-7-3)

$$
P(t) =
$$
\n
$$
\sum_{i=1,3} \sum_{j=a,b} \frac{C_i A}{2} e^{\frac{\sigma^2}{2\tau_i^2} - \frac{t}{\tau_i}} \left[1 + \text{erf}\left(\frac{t - \frac{\sigma^2}{\tau_i}}{\sigma\sqrt{2}}\right) \right] +
$$
\n
$$
\frac{C_i (1 - A) B_j}{2(\frac{\tau_j}{\tau_i} - 1)} e^{\frac{\sigma^2}{2\tau_j^2} - \frac{t}{\tau_j}} \left[1 + \text{erf}\left(\frac{t - \frac{\sigma^2}{\tau_j}}{\sigma\sqrt{2}}\right) \right] -
$$
\n
$$
\frac{C_i (1 - A) B_j}{2(\frac{\tau_j}{\tau_i} - 1)} e^{\frac{\sigma^2}{2\tau_i^2} - \frac{t}{\tau_i}} \left[1 + \text{erf}\left(\frac{t - \frac{\sigma^2}{\tau_i}}{\sigma\sqrt{2}}\right) \right].
$$
\n(4.4)

B. Scintillation Pulse Monte Carlo

We developed a Monte Carlo (MC) that uses the analytical pulse shape model as an input and generates channel-level simulated signals. For a given scintillation pulse size, the number of photons that arrive at the top and bottom arrays are drawn randomly from a binomial distribution, using the top/bottom light collection asymmetry measured with the $CH₃T$ calibration data. The number of photons in a single PMT channel is drawn from a binomial distribution with $p = 1/61$. This is a good approximation for S1 light detection by the 61 PMTs in each array for events in the fiducial volume. The areas of each photon are independently drawn from the single-photon pulse area distribution, averaged across all PMTs. Each photon is then randomly assigned a detection time, drawn from the distribution in Eq. [4.4.](#page-8-1) Photon template functions with the appropriate amplitudes and arrival times are added together to construct a simulated signal. Noise is added to this signal by adding sine waves with frequencies and amplitudes given by the measured noise power spectrum and random phases drawn uniformly on the interval $[0, 2\pi]$. The bandwidth of the DAQ is therefore included in both the noise and the signal by using data-driven signal and noise distributions. The simulated waveform is then sampled at 100 MHz, with the starting point given by a uniform random number between 0 and 10 ns to simulate timing jitter due to the digitizer's sampling. This MC is used in the following sections to simulate scintillation pulses for error analyses and discrimination calculations.

V. ER/NR DATA ANALYSIS AND RESULTS

A. Photon Time Spectrum

We study scintillation characteristics in ER/NR calibration data by constructing photon time spectra that are averaged over many events. Event selection is based on the dark matter search analyses: we study only singlescatter events. These are defined as events with a single

FIG. 6: Average photon detection time spectra for events with pulse area between 40–50 phd $(∼ 11–13 keV_{ee})$. The probability distribution of NR events, from DD neutron data, is shown by the red diamonds and the MC simulated NR time spectrum is shown in red. The measured distributions for ER events from tritium and ¹⁴C are shown in green and blue, respectively. The ER time spectrum generated from MC simulations is shown in blue. The time at which the pulse area reaches 5% of the total, denoted T05 in the text, is used as $t = 0$. The vertical uncertainties in each bin are calculated from Poisson statistics, while the horizontal error bars represent bin width.

S2, preceded by a single S1 within the maximum drift time $(330 \,\mu s)$. To reduce the position dependence of optical transport, indicated by Fig. [5,](#page-7-2) we select only events in a rectangular prism around the beam path of the neutron calibration source, at a median depth of ∼7.5 cm below the liquid surface. The average electric field in this region, calculated using COMSOL electrostatics simulation software [\[43\]](#page-16-29), is 410 V/cm . The times of detected photons in these events are corrected using the channel-to-channel time calibration and the direct-path travel time from the event site to the PMT is subtracted, and weighted by the w_i 's. To align time spectra from different events, we define a common reference of the sample at which the summed waveform crosses 5% of the total pulse area (hereafter denoted T05). The photon time distribution for many events are used to produce average time spectra. As our analysis is focused on the true number of photons arriving in a given pulse, we use S1 pulse areas that are not corrected for position dependent effects in the LUX detector unless otherwise noted.

Average time spectra for the three calibration sources with pulse area between 40–50 phd are shown by the data points in Fig. [6.](#page-8-2) The two electron recoil sources $(^{14}C$ and tritium) show identical spectra, while the nuclear recoil source (DD neutrons) has a spectrum skewed to earlier times. This difference is explored in the context of our analytical model in Section [V B,](#page-9-0) and is used for ER/NR discrimination as discussed in more detail in Section [VI A.](#page-11-0)

The field- and position-dependence of the pulse shapes were studied using the ER calibration data. The electric fields within the detector volume changed significantly between WS2013 and WS2014-16: in the former, the drift field was ∼180 V/cm throughout the detector, while in the latter, the drift field is highly non-uniform, varying from an average of ∼50 V/cm near the bottom to ∼400 V/cm near the top of the detector. For a fixed volume and fixed pulse area, we do not observe a significant difference in the average time spectra for any of the tritium calibrations from WS2013 or WS2014-16. We therefore conclude that there is no significant fielddependence within the limits of our sensitivity. We observe a depth-dependence in the average time spectra, consistent with the expectations from the optical transport model. Since photons from S1 pulses are preferentially detected in the bottom PMT array due to reflections at the liquid-gas interface, we attribute depthdependence of the pulse shapes to the depth-dependence of the geometric efficiency of the bottom array. The effects of the depth-dependence on ER/NR discrimination is discussed in Section [VI A.](#page-11-0)

B. Fits to analytical model

The analytical model was fitted to the average time spectra in both ER and NR data to extract scintillation characteristics. To measure energy dependence, we fit the average time spectra binned by the reconstructed ERequivalent energy (given in keV_{ee}),

$$
E_{rec} = W\left(\frac{S1}{g_1} + \frac{S2}{g_2}\right). \tag{5.1}
$$

Here, S1 and S2 denote the S1 and S2 pulse areas corrected for all position-dependent effects using the ⁸³Kr calibrations. Over the period of the experiment, the parameters g_1 varied between 0.100 ± 0.002 and 0.097 ± 0.001 phd/photon, and g_2 varied between 18.92 ± 0.82 and 19.72 ± 2.39 phd/electron. These are the detector-specific gains for S1 signals and S2 signals during the calibration campaigns used in this analysis [\[1\]](#page-15-0). $W = 13.7$ eV is the average energy required to create either a scintillation photon or an ionization electron in liquid xenon [\[44\]](#page-16-30). For ER events, we use bins of 4 keV $_{ee}$ in E_{rec} from 5–45 keV_{ee}. For NR events, we use bins of 2 keV_{ee} from 5–17 keV_{ee}. To obtain the true energy of events contributing to each E_{rec} bin, the distribution of recoil energies for each source is simulated using the NEST light and charge yield models tuned to LUX data, given in Ref. [\[33\]](#page-16-19) (for ER) and Ref. [\[32\]](#page-16-18) (for NR). We report the mean and the ± 1 σ of the simulated distributions as the true energy and its error. This analysis

includes ERs with true energies from ∼5–46 keV and NRs with true energies from \sim 25–74 keV.

In addition to the timing effects built into the analytical model, the photon time spectra experience a spread due to statistical fluctuations in T05. These depend on the scintillation emission distribution and the total number of detected photons, and produce an additional smearing that could be mistaken for an energy dependence in the underlying time spectra. We model this effect using the MC to simulate events in each energy bin. The distribution of T05 in the appropriate energy bin is convolved with the analytical model before fitting to the measured average time spectra.

Several of the fit parameters are expected to remain constant across energies and particle types. The singlet time constant τ_1 and the Gaussian fluctuation parameter σ are expected to be the same across all energy bins and for both ER and NR data, as they are dependent solely on the scintillation physics of the Xe_2^* dimer and the timing resolution, respectively. Similarly, the optical transport parameters $(A, B_a, \tau_a, \text{ and } \tau_b)$ depend solely on photon transport in LUX, and should be constant across all spectra for a fixed position inside the detector. The values used in these fits are given in Table [I](#page-7-1) for the DD beam location, and are constant for both ER and NR spectra across all energies.

In contrast, the ratio of C_1/C_3 is allowed to vary independently for each energy bin in both ER and NR data. This allows our model to capture the difference in the singlet/triplet ratio between ER and NR events, as well as any possible dependence on recoil energy. We also allow τ_3 to vary between ER and NR datasets to allow it to capture any small recombination effects, as discussed in Section [IV.](#page-6-3)

In order to fit all of these parameters with the appropriate constraints and correlations, we simultaneously fit the average time spectra at all energies using a global log-likelihood given by

$$
\log(L) = \left[\sum_{\text{ER, NR}} \sum_{i=1}^{N} \sum_{j=1}^{M} -\frac{(Y_j - Y_{calc})^2}{\sqrt{2} Y_j} \right], \quad (5.2)
$$

where N is the number of bins in energy, M is the number of bins in the average pulse timing distribution for each energy bin, Y_i is the number of photons in bin j of the timing distribution, and Y_{calc} is the height of bin j, calculated by the model. The singlet and triplet times are constrained to vary within [0 ns, 10 ns] and [18 ns, 35 ns], respectively, to avoid degeneracy. We maximize $log(L)$ using the Minuit optimizer class provided by the ROOT framework [\[45\]](#page-16-31).

The results of the global fit are summarized in Table [II.](#page-10-0) The statistical errors are those returned by the fit routine. In addition, there are two main sources of systematic error: errors introduced in the analysis and fit procedure, and errors from uncertainties in the optical model used to produce A, B_a , τ_a , and τ_b . We quantify the first by performing the analysis on simulated events

TABLE II: Summary of parameters used in fitting our photon time spectra. Expected values in column 2 come from the literature where appropriate: the predictions for the triplet and singlet times come from the average values in Ref. [\[12\]](#page-16-0), while the prediction for $C_1\tau_1/C_3\tau_3$ comes from the measurement in Ref. [\[11\]](#page-15-9). The expected value of σ is calculated from Eq. [4.3.](#page-7-3) The best-fit parameters are shown with $\pm 1 \sigma$ statistical uncertainties, calculated from the fit. The analysis systematic is computed by performing the same fit procedure on simulated data with known input parameters. The optical transport systematic uncertainty comes from varying the optical model used in the fit, i.e. fitting A, B_a , τ_a , and τ_b to optical simulations with the $\pm 1 \sigma$ extremes on optical parameters from Ref. [\[38\]](#page-16-24).

| Parameter | Expected | Fit constraint | Best fit \pm stat. | Analysis sys. | Optical transport sys. |
|---------------------------|--------------------|----------------|----------------------|------------------|---------------------------|
| $(C_1\tau_1)/(C_3\tau_3)$ | ~ 0.1 (ER) | none | 0.042 ± 0.006 | $\pm 3.1\%$ | $+75%$ -66% |
| | none (NR) | none | 0.269 ± 0.034 | $\pm 3.1\%$ | $+20%$ -10% |
| τ_1 | 3.1 ± 0.7 ns | $0-10$ ns | 3.27 ± 0.66 ns | $\pm 1.0\%$ | $+11%$ $-70%$ |
| τ_3 | 24 ± 1 ns (ER) | $18-35$ ns | 25.89 ± 0.06 ns | $\pm 1.9\%$ | $+0.5\%$ -0.6% |
| | 24 ± 1 ns (NR) | $18-35$ ns | 23.97 ± 0.17 ns | $\pm 1.9\%$ | $+0.1\%$ -1.1% |
| А | 0.0574 | fixed | | | $+34%$ -9.1% |
| \boldsymbol{B} | 1.062 | fixed | | | $+1.9\%$ -1.0% |
| τ_a | 11.1 ns | fixed | | | $+17%$ $-21%$ |
| τ_b | 2.70 ns | fixed | | | $+0\%$ -8% |
| σ | 3.2 ns | none | 3.84 ± 0.09 ns | $\pm 1.1\%$ | $+1.2\%$ -1.2% |
| | | | | | |

with known input parameters. Shifts between the input parameters and the reconstructed parameters are quoted as the systematic errors shown in the fifth column in Ta-ble [II,](#page-10-0) and are $O(2\%)$. The errors due to uncertainties in the optical model is quantified using the χ^2 distribution of the optical model fit given in Ref. [\[38\]](#page-16-24). Of the seven free parameters in the optical model, we find that only two affect the photon transport times: the liquid xenon absorption length and the teflon reflectivity. We run new optical transport simulations allowing these parameters to vary along the $\Delta \chi^2 = 8.18 \,(1\sigma)$ contour of the optical model parameter space. We then propagate these new simulated distributions into our pulse shape model and redo the analysis to extract new pulse shape parameters. We report the variations from the best-fit values as the systematic error, which is listed in the sixth column of Table [II.](#page-10-0) This is the dominant error in our analysis.

The best-fit singlet/triplet ratios as a function of energy are shown in Fig. [7.](#page-11-1) For electron recoils we find $(\tilde{C_1}\tau_1)/(C_3\tau_3) = 0.042 \pm 0.006$ (stat) $^{+0.092}_{-0.034}$ (sys), averaged across all measured energies, which is lower than existing results in the literature. We note this is the first measurement of the singlet/triplet ratio with both a low energy ER source and an applied electric field. The energy dependence at zero field, measured by the XMASS collaboration (shown by cyan diamonds in Fig. [7\)](#page-11-1) is correlated with a lengthening of the long time constant from 28 ns to 32 ns, which suggests that they are observing an increase in the recombination-related time constant with energy which are not explicitly accounted for in their model. We do not observe an energy dependence in neither the time constant nor in the singlet/triplet ratio, consistent with

the hypothesis that the applied electric field in our experiment suppresses recombination contributions to the pulse shape. For nuclear recoils, we find $(C_1\tau_1)/(C_3\tau_3)$ $= 0.269 \pm 0.034 \text{ (stat)} \, \frac{+0.182}{-0.083} \text{ (sys)}$, averaged across all energies probed in this analysis. The only analogous measurement in the literature uses recoiling fission fragments and finds $(C_1 \tau_1)/(C_3 \tau_3) = 1.6 \pm 0.02$, though in a vastly different energy regime at O(100 MeV) [\[7\]](#page-15-5). Our result is therefore the first nuclear recoil singlet/triplet ratio measurement that is directly relevant for dark matter TPC experiments.

We test for energy dependence of the singlet/triplet ratio by fitting both a constant value and a power law dependence, the latter given by $(C_1 \tau_1)/(C_3 \tau_3) = \alpha E^{\beta}$. Such an energy dependence is well-established in liquid argon [\[46\]](#page-16-32), but has never been directly explored in xenon. For electron recoils, the best-fit values of the power law give $\alpha = 0.063$ and $\beta = -0.12$. The $\chi^2/\text{d.o.f.}$ for the constant and power law models are $16.6/9$ (p = 0.06) and $13.7/8$ ($p = 0.09$), respectively. For nuclear recoils, the best-fit values of the power law give $\alpha = 0.15$ and $\beta = 0.15$. In this case, the $\chi^2/\text{d.o.f.}$ for the constant and power law models are $4.6/5$ (p = 0.47) and $3.2/4$ $(p = 0.52)$. We conclude that our data is statistically consistent with both models, and both are compared to data in Fig. [7,](#page-11-1) Fig. [9,](#page-13-0) and Fig. [10](#page-14-0) for completeness.

Our best-fits of the triplet and singlet time constants, τ_1 and τ_3 , agree with previously measured values. The expected values, listed in Table [II,](#page-10-0) are the error-weighted averages computed in Ref. [\[12\]](#page-16-0) based on a survey of measurements in the literature. The only value in slight tension is the triplet time constant that we measure for electron recoils, which is higher than both the expected value and our best-fit for nuclear recoils. This is consistent with small recombination effects that are not accounted for in our model. If we assume $\tau_{3_{er}} = \tau_{3_{nr}} = 23.97$ ns and take the recombination time distribution derived in Ref. [\[8\]](#page-15-6) $(P(t) \propto [1 + (t/\tau_R)]^{-2})$, simulations reproduce our bestfit distribution with $\tau_R \approx 0.6$ ns. This expression for recombination time may not be directly applicable here, as it is derived by solving a diffusion equation with no applied electric field. However, we note that the qualitative agreement with the empirical prediction of $\tau_R = 0.7$ ns from Ref. [\[12\]](#page-16-0) is encouraging. Regardless, our result for $\tau_{3_{er}}$ is still within the range of τ_3 measurements available in the literature [\[6,](#page-15-4) [11\]](#page-15-9), indicating that recombination plays a minor role in the pulse shapes for electron recoils in our experiment.

VI. PULSE SHAPE DISCRIMINATION

A. Prompt Fraction Discriminator

To discriminate between the two types of events we adapt a Prompt Fraction Discriminator (PFD), a standard technique which has been successfully adapted for use in other liquid xenon and liquid argon dark matter experiments [\[11,](#page-15-9) [16,](#page-16-33) [17,](#page-16-34) [47,](#page-16-35) [48\]](#page-16-36). The parameter is defined as:

$$
PF = \frac{\int_{t0}^{t1} S1(t)dt}{\int_{t2}^{t3} S1(t)dt} = \frac{\sum \text{Prompt Photons}}{\sum \text{Total Photons}}.
$$
 (6.1)

The four variables, $t0$, $t1$, $t2$, and $t3$, are allowed to vary independently in the range of −30 to 170 ns to minimize the leakage of ER events into the NR acceptance region, where the ∼50% NR acceptance region is defined as everything above the NR median (\overline{NR}) . No additional constraints on these parameters were imposed and cases where $t0 > t2$, etc. were explored. For each combination of times, the PF is applied to the calibration data and we create a map of the fraction of event that appear in the NR acceptance region. We choose the combination of times that produces the minimum leakage as our PF and the NR as our optimal Prompt Fraction Discriminator (PFD).

To calculate an unbiased performance of the PFD we separate the calibration datasets into two groups and apply a weighting to mitigate the energy dependence of the source. The calibration datasets are divided into 10 phd bins before events in each bin are randomly assigned to either the training or the testing group. Both groups contain 50% of the data and there is no statistically significant difference between their average detected photon time spectra, position, or energy distributions. The training group of events are used to train our PFD. The results of this PFD are applied to the events in the testing group and are presented in this paper. Since the count

FIG. 7: Singlet/triplet ratio $(C_1\tau_1/C_3\tau_3)$ measured for nuclear recoils (Top) and electron recoils (Bottom) using LUX calibration data. Only statistical uncertainties are shown. Calibration sources are DD neutrons (red), tritium (blue), and ${}^{14}C$ (green). Measurements in different energy bins are shown by the square points, while the average across all energies is shown by the solid line. A power law is fit to the data is presented by the dashed line. We also show measurements of the ER singlet/triplet ratio at zero field from Ref. [\[11\]](#page-15-9) (cyan diamonds), and a measurement using a 207 Bi internal conversion source at 4 kV/cm from Ref. [\[8\]](#page-15-6) (purple diamond). In Ref. [\[11\]](#page-15-9), the singlet fraction (denoted F_1) is given rather than the singlet/triplet ratio. For direct comparison to this work we make the conversion $(C_1\tau_1)/(C_3\tau_3) = F_1/(1+F_1).$

rate of the calibration sources is energy dependent, each 10 phd bin is weighted equally when calculating the total leakage and is not weighted by the number of events in that particular bin. We optimize the PFD for a 'flat' distribution in pulse area.

Using the six different calibration campaigns carried out at various times during WS2013 and WS2014-16, an optimal PFD has been found. ¹⁴C was only injected at the end of the LUX WS, whereas tritium and DD calibrations were carried out during each of the five calibrations campaigns. Carrying out the PFD optimization using

FIG. 8: Normalized prompt fraction distributions for NR (red) and ER (blue) events with raw S1 pulse areas from 40–50 phd. Solid red and blue traces are the simulated NR and ER prompt fraction distributions for this pulse area. The vertical error bars indicate the Poisson uncertainty, while the horizontal error bars represent bin width. The solid black line indicates the median of the NR distribution, and the prompt fraction region above the DD median is considered to be the NR acceptance region. $26 \pm 2\%$ of the ER distribution was found to lie within this region.

the events in the training group, which is a composite of each of these calibration campaigns, gives an optimal prompt window of −8 to 32 ns and total window of −14 to 134 ns. For this optimal combination of prompt and total photons, an example of PF values in the 40–50 phd bin applied to events the testing group is presented in Fig. [8.](#page-12-0) This PFD is also applied to events generated using the MC simulation and shows consistency with data. When this PFD is trained on the individual campaigns the optimal windows are found to vary up to 6 ns and are considered statistical fluctuations rather than real changes in the photon detection time spectrum.

Figure [9](#page-13-0) shows the PF distributions calculated from DD, CH3T and ¹⁴C calibration data and events generated using the MC simulations. At small pulse areas, the PF for both experimental and simulated data have a large spread due to the low photon statistics in each event. At larger pulse areas, more photons are reconstructed and the photon time spectra for individual events will appear more ER or NR-like. This results in a narrowing of the bands for the two populations at higher energies. This reduction of band width provides a lower leakage fraction at higher energies.

Fig. [10](#page-14-0) shows the fraction of ER events that leak into the NR acceptance region. The DD calibrations were used to compute the 50% NR acceptance region while the ¹⁴C and CH₃T calibrations were used to calculate the leakage into this region. For the distributions shown, the leakage in the lowest energy bin of 10–20 phd is $39.4 \pm 2.7\%$. In the 40-50 phd bin, the highest bin

used in the WS2013 and WS2014-16 analysis, the leakage fraction is reduced to $26.1 \pm 2.0\%$. The leakage fraction continues to decrease at higher energies.

Using calibrations, we study the vertical position dependence on our ability to discriminate between ER and NR events. At greater depths in the detector the PFs move to larger values as more photons are detected at the PMTs with less scattering. This geometric affect applies to both the ER and NR events and causes both bands to move by a similar value in PF for a given number of detected photons. Thus, we do not observe any depth dependence in the leakage fraction. The leakage fractions at different depths in the detector are consistent with those presented in Fig. [10.](#page-14-0) The overall leakage fraction, for a flat distribution of events up to 200 phd, generated from simulations, is 22.9%.

B. Two Parameter Discrimination for Dark Matter Searches

To improve the ability of LUX to discriminate ER from NR events, the PFD can be used in conjunction with the charge-to-light ratio $(\log_{10}(S2/S1))$. This is shown in Fig. [11](#page-14-1) for events with pulse areas between 40–50 phd. An elliptical region, centered at the median values of the NR distribution, is chosen to include ∼90% of the total distribution. In this two dimensional space, a line passing through the median of the NR population is used to create a linear cut to discriminate between the two populations of events. The region, above this linear cut, away from the ER population, is defined as the NR acceptance region. In this manner, ∼50 % of the NR acceptance region is preserved by the ellipse and the vertical cut.

For each 10 phd bin, the three free parameters of the ellipse (inclination and two radii) and the one free parameter of the linear cut $(x \text{ or } y\text{-intercept})$ are varied to minimize the ER leakage into the NR acceptance region. To test leakage, a line is drawn through the center of the ellipse with either a x-intercept between 1.25 and 2.75 or a y-intercept between 0 and 1. In this way, a cut that closely resembles just the charge-to-light discrimination is also tested. When this method is applied to 10 phd wide bins between 10–100 phd, the ellipse's inclination and x-intercept with the least leakage are often very similar. For each of these bins, the optimal inclination of the NR population and x-intercept and are close to 30° counter-clockwise w.r.t. the vertical. and 1.7, respectively. As the pulse area of the event gets smaller, the distribution of possible PF and charge-to-light increase, and thus the optimal radii for the ellipses vary to capture this change.

For the example shown in Fig. [11](#page-14-1) for 40–50 phd, the overall ER leakage is reduced by 50 % compared to just the $\log_{10}(S2/S1)$. Using the $\log_{10}(S2/S1)$ bands for this population of events [\[1\]](#page-15-0), an overall ER leakage of 0.4 ± 0.1 % is achieved. When the $\log_{10}(S2/S1)$ and

(a) NR data and the constant value model.

(b) NR data and the power law model.

(c) ER data and the constant value model.

(d) ER data and the power law model.

FIG. 9: PF distribution for NR events (Top), and ER events (Bottom). The red and blue dots indicate the NR and ER bands respectively (median and median ± 34 percentile) calculated from data. The traces indicate the bands calculated from simulated data using the constant value model (Left) and the power law model (Right). The solid traces indicate the medians and the dashed lines indicate the median ± 34 percentile. We define the region above the \overline{NR} as the NR acceptance region.

PFD are combined to produce a new discrimination parameter, as described above, the overall ER leakage reduces to 0.3 ± 0.1 %. The comparison of the various discrimination methods are presented in Table [III.](#page-14-2)

VII. CONCLUSION

We have described an analysis of liquid xenon scintillation pulse shapes and the discrimination power in the LUX dark matter experiment. We have developed software that allows for the precise reconstruction of photon detection times within a pulse with an accuracy of 3.8 ns. LUX calibration data from DD, CH_3T , and ¹⁴C sources are used to characterize the photon detection time spectra for NR and ER events at various depths in LUX. Average time spectra are fitted with an analytical model to extract singlet-to-triplet ratios and singlet and triplet decay times. It is found that the singlet-to-triplet ratio for ER events is consistent with the literature within errors. We have made a first measurement of the NR singlet-totriplet ratio at low energies and a non-zero applied electric field. Different τ_3 time constants are found for ER and NR events. We interpret this as residual recombination timing effects, which are not included in our model, adding a small smearing to the ER pulse shape which gets captured by the τ_3 parameter in our fits. These measurements and the reconstructed physical properties of xenon scintillation are relevant for liquid xenon dark matter search experiments, and can inform simulation packages, such as NEST, that are used by the community to compute event distributions in current and future experiments.

TABLE III: Summary of ER leakage percentage into the NR acceptance region using different methods. The data presented here are from all the DD, $\overline{CH_3T}$ and $\overline{14}C$ calibrations from the volume around the DD beam. In all cases the leakage is defined as the number of ER events that occur within the NR acceptance region. The errors indicated are from Poisson statistics. Column 2 presents the charge-to-light yield discriminator from WS2014-16, which is not studied above 50 phd. Column 4 presents performance of the PFD calculated from simulations using a linear singlet-to-triplet ratio.

| S1 Pulse Area [phd] | $Log_{10}(S2/S1)$ [%] | PFD Data $[\%]$ | PFD Simulation $[\%]$ | Two Parameter Data [%] |
|------------------------|-----------------------|---------------------------|---------------------------------|---------------------------|
| $10-20$ | 0.5 ± 0.2 | 39.3 ± 2.7 | 32.7 | 0.4 ± 0.2 |
| 20-30 | 0.4 ± 0.2 | 31.3 ± 2.2 | 29.4 | 0.3 ± 0.1 |
| $30 - 40$ | 0.4 ± 0.2 | 28.9 ± 2.2 | 26.9 | 0.2 ± 0.1 |
| $40 - 50$ | 0.3 ± 0.2 | 25.6 ± 2.0 | 24.5 | 0.1 ± 0.1 |
| $50 - 60$ | | 22.7 ± 2.0 | 22.9 | 0.1 ± 0.1 |
| 60-70 | | 21.7 ± 1.9 | 21.3 | 0.0 ± 0.1 |
| 70-80 | | 19.2 ± 1.8 | 20.2 | 0.0 ± 0.1 |
| 80-90 | | 20.1 ± 1.7 | 19.3 | 0.1 ± 0.1 |
| 90-100 | | 17.9 ± 1.5 | 18.5 | 0.0 ± 0.1 |

FIG. 10: Fraction of ER events that leak into the NR region. The points show the leakage fraction obtained from the data while the solid trace indicates the leakage fraction calculated from simulation. The NR is used to define the ∼50% NR acceptance region. The horizontal error bars indicate the bin widths. The counting errors from Poisson statistics and the errors in \overline{NR} are added in quadrature to calculate the error in leakage fraction. For both data and simulations, the same PFD windows are used.

The photon timing module is applied to calibration data to construct photon detection time spectra. The difference between ER and NR time spectra is exploited to formulate a ratio of prompt to total photons to discriminate ER and NR events. This discrimination pa-

FIG. 11: Two parameter event discrimination using PFD and $log_{10}(S2/S1)$. The distribution contains NR (red) and ER (blue) events with raw S1 pulse areas between 40–50 phd. The yellow ellipse indicates the 90% NR region and the solid black line indicates the discriminator. The region to the left of the black line is defined as the NR acceptance region and is chosen to preserve 50% of the ellipse's NR region while minimizing the number of ER events in this region.

rameter (PF) is optimized, using a training data set, to minimize the leakage of ER events into the ∼50% NR acceptance region. The photon detection time and the prompt fraction distributions are shown to agree with those generated from the MC simulations using the bestfit analytical model, allowing us to extrapolate to energy regions where no calibration data is available. The discrimination power of the PFD improves with energy. For a flat distribution of events in the WS region from 10– 50 phd, the ER leakage is found to be 31.3 %. Between 10–200 phd, the total leakage is 25.2 % .

In the two-dimensional parameter space composed of charge-to-light ratio $(\log_{10}(S2/S1))$ and PF, an improved discriminator is developed. This discriminator is required to preserve ∼50 % NR acceptance while reducing the ER leakage into the region. Over the WS region of 10–50 phd, the ER leakage, measured using the charge-to-light discriminator, is 0.4 ± 0.1 % and reduces to 0.3 ± 0.1 %, measured using the two parameter discriminator. This discrimination power increases at higher energies, making it an attractive background reduction technique for dark matter searches looking for nuclear recoils at higher energies than the traditional WIMP search. Examples of these dark matter searches include models in which dark matter scatters inelastically, or with a momentumdependent cross-section. Using the parameters from the analytical model, the pulse shape discrimination bands can be extrapolated out to higher energies than accessible by calibration sources, or can be extrapolated to assess the pulse shape discrimination power in future liquid xenon experiments.

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