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ARISE: An Algorithm for Rapid Ion Spectrum Extraction Enabling Real-Time Optimisation in High-Repetition-Rate Laser-Driven Ion Acceleration

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Abstract

Recent advances in high-power, high-repetition-rate laser systems are driving the adoption of data-driven experimental approaches in high-energy density science. To fully realise the potential of these methodologies, automated and high-throughput analysis of key diagnostics is essential for effective feedback and real-time optimisation. We present a novel algorithm, ARISE (Algorithm for Rapid Ion Spectrum Extraction), developed for fast and reliable extraction of laser-accelerated ion spectra from Thomson parabola spectrometers, capable of operating at repetition rates exceeding 20 Hz. ARISE enables real-time, data-driven experimentation through features including background subtraction, automatic identification of the zero-deflection reference point, and automated determination of maximum ion energy. We validate the accuracy of ARISE in spectrum extraction and energy detection, and demonstrate its integration within a Bayesian optimisation framework during a proof-of-concept experiment conducted using the 350 TW SCAPA laser, enabling real-time optimisation of laser-accelerated ion beam parameters.

1. Introduction

Beams of high-energy ions can be generated during intense laser-solid interactions, with maximum proton energies exceeding 100 MeV demonstrated to date[1–3]. Novel features such as ultra-short bunch duration mean these ion sources have a wide array of potential applications, including fast-ignition inertial confinement fusion (ICF)[4,5], proton radiography[6], radiation damage testing[7,8], and radiobiological research[9–11]. Until now, much of the development in laser-driven ion acceleration has been undertaken using high-energy (tens-to-hundreds of Joules) lasers and targetry systems that were inherently limited to low ($\ll$1 Hz) repetition rates[12–14]. This limitation has significantly hindered the implementation of statistical and data-driven methodologies, which could be crucial for optimising and stabilising ion acceleration mechanisms.

Supported by advances in high-power laser technology in recent years, a number of petawatt-class high-repetition rate laser systems have been developed that are capable of $\geq$1 Hz operation[15]. Such systems have motivated the development of rapidly replaceable targetry such as tape-drive systems[16], liquid crystals[17], cryogenic[18] and liquid jets[19]. As a consequence, it is now possible to produce high-energy laser-driven ion beams at 1 Hz, and beyond (see for example reference[20]). These efforts have enabled statistical and data-driven experimentation, including machine learning (ML)-based optimisation of the ion source[21,22] and the generation of training data sets for neural networks to build surrogate models[23–26]. These emerging approaches will support the active optimisation and stabilisation of laser-plasma ion sources, facilitate deeper investigation of the underpinning physics and enhance experimental design - as demonstrated in inertial confinement fusion (ICF) experiments[27]. To fully realise data-driven ion acceleration experiments, it is essential to automate the extraction of key ion beam parameters — such as the energy spectrum, conversion efficiency and maximum energy — at a rate that matches or exceeds the laser repetition rate.

In this article, we present the *algorithm for rapid ion spectrum extraction* (ARISE) - a software tool designed for extracting ion energy spectra from a Thomson parabola spectrometer at Hertz-scale repetition rates, enabling data-driven experimentation with live-feedback. ARISE incorporates background subtraction, automatic identification of the zero-deflection reference point (defining the origin of parabolic ion tracks), and automatic determination of the maximum ion energy. At the Scottish Centre for the Application of

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Plasma-based Accelerators (SCAPA) at the University of Strathclyde, we have developed a laser-driven ion acceleration beamline, which uses the 5 Hz, 350 TW laser for investigation of the physics underpinning laser-driven ion acceleration and for applications such as ultra-high dose rate radiobiology[28] and radiation damage studies[29,30]. We demonstrate the performance of the ARISE algorithm using the SCAPA laser–ion beamline operating at 0.2 Hz, a rate constrained solely by data transfer speeds and the readout time of the diagnostic camera. Integrated into a feedback loop, ARISE performed real-time ion spectral analysis to autonomously guide and optimise the maximum energy of laser-accelerated protons. Furthermore, we show that ARISE can process ion spectra at rates exceeding 20 Hz when applied to representative archival experimental data. We validate its accuracy in both spectrum extraction and automatic detection of the maximum proton energy ($E_{\mathrm{p,max}}$), highlighting its suitability for high-throughput ion diagnosis in data-driven laser–plasma experiments.

2. Ion Spectrometer Design and Data Capture

A Thomson parabola spectrometer (TPS) separates ions based on their kinetic energy and charge-to-mass ratio, producing characteristic parabolic traces[31]. In this work, we employed the TPS design reported in Carroll *et al.*[32], the operating principle of which is illustrated in figure 1. The spectrometer featured a 500 μm diameter pinhole to limit angular acceptance. Ions were deflected by a 0.6 T magnetic dipole pair and a pair of wedged electric plates, across which a total potential difference of 5 kV was applied. The ions were then detected by a microchannel plate (MCP) coupled to a phosphor screen (Hamamatsu F2226-14PF143). The end of the electric plates was positioned 75 mm upstream of the MCP. The resulting parabolic ion traces were recorded using a 16-bit sCMOS camera (Andor Neo) and example measurements are shown in figure 2(a). The geometry and dimensions of the magnetic and electric field regions matched those described in Carroll *et al.*[32].

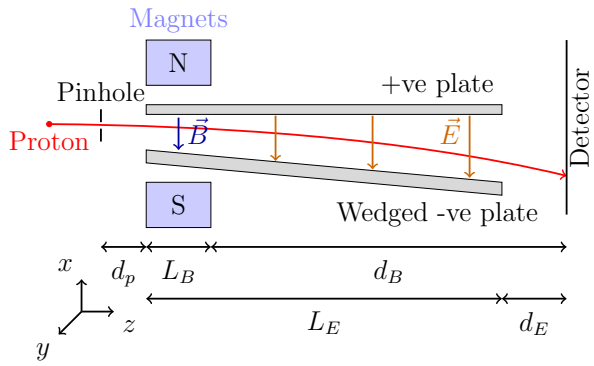

Figure 1 : Schematic view of the Thomson parabola spectrometer in the x-z plane. The electric and magnetic fields are both oriented along the x-axis, resulting in ion dispersion along the x- and y-axes due to their respective influences.

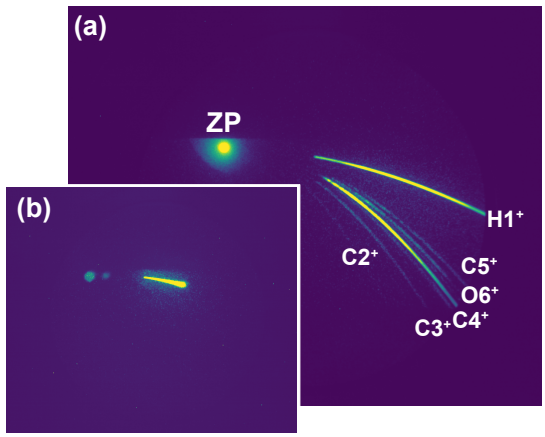

Figure 2 : (a) Representative MCP phosphor screen image showing the reference zero point (ZP), corresponding to undeflected neutral atoms and X-rays, and a set of ion tracks (H - hydrogen, C - carbon, O - oxygen, with given charge states). (b) An example MCP phosphor screen image showing the filtered ion signal through 200 μm-thick Mylar, for the energy calibration.

2.1. *Charged Particle Trajectories in a Thomson Parabola Spectrometer*

Charged particles entering a Thomson parabola spectrometer are deflected by electric and magnetic fields, with their motion governed by the non-relativistic Lorentz force:

$$\vec{F} = q\left(\vec{E} + \vec{v} \times \vec{B}\right) \tag{1}$$

where q is the particle charge and $\vec{v}$ is its velocity, $\vec{E}$ and $\vec{B}$ are the electric and magnetic fields, respectively.

In the small-angle approximation, the transverse displacements resulting from the electric and magnetic fields — accounting for a drift region between the field termination and the detector — are approximately given by:

$$x \sim \frac{q|\vec{E}|L_E}{mv_z^2}\left(\frac{L_E}{2} + d_E\right) \tag{2}$$

$$y \sim \frac{q|\vec{B}|L_B}{mv_z}\left(\frac{L_B}{2} + d_B\right) \tag{3}$$

where L_E and L_B are the effective lengths of the electric and magnetic field regions, d_E and d_B are the distances from the end of each field region to the detector plane, m is the mass of the particle, and v_z is the longitudinal component of the velocity.

These expressions describe the characteristic parabolic traces observed in a TPS under the assumption of uniform fields. However, in the present setup employing

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a wedged electrode design, the fields are inherently non-uniform, necessitating a more general approach to modelling ion trajectories.

Assuming a constant particle mass m, and defining the velocity and position vectors as $\vec{v} = (v_x, v_y, v_z)^T$ and $\vec{r} = (x, y, z)^T$, Newton's second law provides the time evolution of the velocity:

$$\frac{d\vec{v}}{dt} = \frac{q}{m}\left(\vec{E} + \vec{v} \times \vec{B}\right) \tag{4}$$

To determine the full trajectory, we also evolve the position:

$$\frac{d\vec{r}}{dt} = \vec{v} \tag{5}$$

Rather than integrating these equations in time and checking when the particle reaches a fixed longitudinal position (e.g., $z = z_{\text{det}}$), it is often advantageous to reparameterise the system using z as the independent variable. This approach simplifies numerical integration in systems where z increases monotonically.

Applying the chain rule:

$$\frac{d}{dt} = \frac{dz}{dt} \cdot \frac{d}{dz} = v_z \frac{d}{dz} \tag{6}$$

the equations of motion can be rewritten as:

$$\frac{d\vec{r}}{dz} = \frac{\vec{v}}{v_z} \tag{7}$$

$$\frac{d\vec{v}}{dz} = \frac{q}{m v_z}\left(\vec{E} + \vec{v} \times \vec{B}\right) \tag{8}$$

Equations (7) and (8) form a coupled system of six first-order ordinary differential equations (ODEs) with respect to z: three governing the components of velocity and three for the components of position.

The equations of motion are then solved numerically using the `solve_ivp` solver from the `scipy.integrate` module in Python[33], over a predefined range of particle energies and charge states for the electric and magnetic fields determined by the TPS design geometry, potential difference and magnetic field strength. The resulting transverse displacements (x, y) at the detector produce the characteristic parabolic traces of a TPS.

2.2. *Data Capture*

To enable data acquisition and preparation for ARISE at the necessary repetition rate, we employ custom data capture and management software. This system facilitates real-time acquisition and handling of all laser data, metadata, and diagnostics associated with a given experiment. It also supports automated control of the laser system and target delivery, and can execute grid scans or implement Bayesian optimisation algorithms[35,36]. When combined with ARISE, this analysis framework enables fully automated, real-time optimisation of key proton beam parameters on the SCAPA beamline, as elaborated in section 4.1.

3. ARISE Structure and Functionality

This section outlines the core functionality provided by ARISE for the automated extraction of ion spectra from unprocessed images of the TPS parabolae. Key features include automatic background subtraction, zero-point detection, spectral extraction, and identification of the maximum ion energy. The implementation of these components for spectral analysis is illustrated schematically in figure 3.

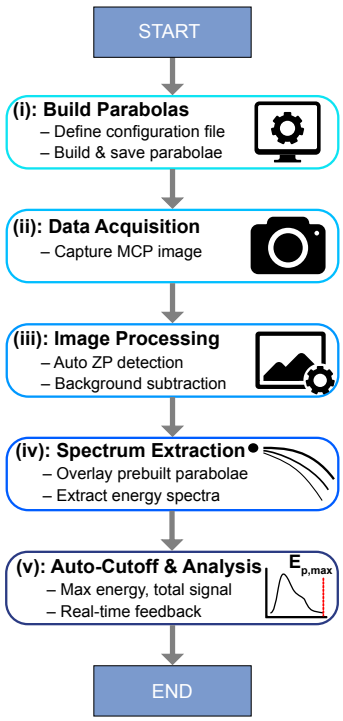

Figure 3 : Flow diagram illustrating the main components of the ARISE processing pipeline. The workflow begins with user-defined parameters specified in the configuration file, followed by data acquisition from the CCD (MCP image). The image then undergoes a series of processing steps, including zero-point detection, cropping, rotation, and background subtraction. The resulting image is subsequently passed to the spectral extraction module and the automated detection of $E_{\text{p,max}}$.

The various steps are as follows:

(i): The initial stage of ARISE analysis requires the user to define fixed design parameters of the TPS within a configuration file. These parameters include the spectrometer geometry, the potential difference across the electrode plates, magnetic field strength, the spatial calibration of the camera, and the range of ion energies to be extracted.

Based on this configuration, ARISE constructs the predicted ion parabolic tracks for a given TPS design, ion species, and specified energy range. For each predefined energy increment, the ion deflection coordinates are computed using a numerical solver for the coupled system of ordinary

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differential equations (ODEs) described in section 2, which models the ion trajectory through the regions L_E, L_B, d_E and d_B. The resulting deflections are first calculated in real space and subsequently converted into pixel coordinates using the spatial calibration. These coordinates define sampling paths along the parabolic tracks, enabling the extraction of ion signal intensity corresponding to discrete energy bins.

One of the inherent limitations of a TPS is the potential overlap of neighbouring species' tracks as they converge at higher ion energies. To address this, the ARISE algorithm issues an alert when track overlap is detected, enabling the user to take corrective action to minimise such occurrences—for example, by increasing dispersion through a higher field strength or by positioning the detector further from the fields. The algorithm includes a dedicated class that can be activated by setting a corresponding flag in the configuration file to *True*, which enables the detection of ion species overlap. A second flag can similarly be set to *True* to construct truncated parabolas; that is, the parabolas are terminated at the point of overlap, beyond which the ion species can no longer be identified with certainty. This truncation feature can be enabled via the configuration file, and the newly generated parabolas can be stored for subsequent use.

Due to the nonlinear relationship between particle energy and deflection, the spatial separation between adjacent energy bins decreases at higher energies. To mitigate the resampling of identical pixels across adjacent bins, ARISE removes duplicate pixel coordinates and assigns each pixel a representative energy value, calculated as the average of its contributing energies.

While ion trajectories (parabolae) can be recalculated for each newly acquired signal, doing so can be computationally intensive depending on the energy resolution, thereby limiting the repetition rate of analysis. However, as the ion deflection paths depend only on fixed geometrical and field parameters — which are assumed to remain constant between shots — ARISE addresses this limitation by storing precomputed parabolae for each TPS configuration. This significantly reduces processing speed by enabling their reuse in subsequent analyses.

(ii): Once the ion trajectories have been either calculated or loaded from a prebuilt parabola file, unprocessed MCP images can be acquired for analysis.

(iii): The acquired image of the MCP phosphor screen is then processed to extract the ion spectra. ARISE includes an automatic detection routine to identify the point of zero deflection — hereafter referred to as the zero-point (ZP) — which typically consists of undeflected neutral particles and X-rays. The ZP represents the origin of the processed image and is the convergence point of all ion parabolae. An example of the ZP position relative to the ion tracks is shown in figure 2(a).

The ZP detection feature initially employs image erosion techniques (see for example[37]) to eliminate saturated pixels and then identifies the coordinates of maximum intensity. Optionally, a region of interest (ROI) can be specified around the expected ZP location, reducing computation time and minimising interference from other bright regions in the image. The approximate ZP coordinates are defined in the configuration file, from which ARISE constructs the ROI for targeted erosion and peak detection.

Once the ZP has been identified, the next stage involves subtracting the background from the image. Initially, a median filter (see for example[33,38]) may be applied across the entire image. This filter computes, for each pixel, the median value within its surrounding region—defined by a user-specified kernel size—in order to suppress impulse noise such as isolated high-signal pixels resulting from detector noise or high-energy X-rays, as well as to mitigate the effects of dead or abnormally responsive pixels, while preserving sharp edges. The filtering process has been implemented in a multi-threaded manner to enhance computational efficiency by dividing the image into smaller segments—whose number is specified in the configuration file—processing them in parallel, and subsequently reconstructing the full image.

Empirical analysis of hundreds of measurements indicated that assuming a radially symmetric background centred on the ZP provides the best agreement with observed background signal characteristics. To ensure accurate background estimation, regions beyond the phosphor screen and areas surrounding the ion tracks are masked out to avoid contamination of the background model. A saturation check is then performed within the ion track region, during which the user is notified if saturation occurs and to what extent. Specifically, if the proportion of saturated pixels exceeds a user-defined threshold (for example, $>1\%$), an alert is issued, prompting the user to adjust the MCP amplification accordingly. Then, for each radial distance from the ZP, the average signal value is calculated and adjusted using a standard deviation multiplier, configurable by the user, to determine the threshold for background subtraction. In our dataset, we found that subtracting values exceeding the average by four standard deviations yielded the most reliable results in terms of isolating the ion signal.

Alternative methods for background subtraction — including Otsu's method[39], sampling from tracks parallel to the ion trajectories, and background region-of-interest sampling — were also investigated. However, these approaches were found to perform less effectively, both in terms of background removal accuracy and computational efficiency.

(iv): Following background subtraction, the ion spectrum is extracted along the previously defined ion tracks. For each track, two additional bounding parabolae are constructed to enclose the full spatial width of the ion signal, determined by the dimensions of the TPS pinhole. The sampled track

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width, δ, is calculated using the TPS geometry as:

$$\delta = d + (s + d)\,\frac{L_E + d_E + d_p}{L} \tag{9}$$

where d is the pinhole diameter, s is the effective source size, d_p is the distance from the pinhole to the start of the electric plates, and L is the source-to-pinhole distance[40]. At each energy step, a lineout is extracted perpendicular to the trajectory defined between the bounding parabolae. The sum of pixel values along this lineout provides a measure of signal intensity as a function of ion energy, corresponding to the relative number of detected ions.

While it is possible to convert these pixel counts into absolute ion numbers using established calibration methods[41–43], an absolute calibration has not yet been implemented in the current version of ARISE. Work is underway to perform such a calibration, following methodologies similar to those described in Harres *et al.*[43].

(v): Finally, from the extracted spectrum, key physical quantities are derived in real time, including the maximum ion energy and total signal intensity, corresponding to the total ion flux. The approach can easily be extended to extract other quantities, such as the spectral temperature of the distribution. These outputs are used both to inform experimental decision-making and to support data-driven optimisation routines aimed at enhancing the quality and performance of the ion source.

4. Code Performance and Validation

In this section, we evaluate the performance of key features within the ARISE framework, through both verification tests and application to real-time experimental data.

Verification of model accuracy and energy measurements: A central requirement of ARISE is the robust and reliable automatic detection of the maximum proton energy for a given proton spectrum. This capability is critical for supporting optimisation and stability studies of laser-driven ion sources. Prior to implementing and assessing automated detection routines, we first validated the accuracy of the ion energies predicted by the ARISE deflection model.

To validate both the accuracy of the ion deflection model within ARISE and the energy resolution of the TPS, selected regions of the MCP were covered with Mylar foils of known thickness, designed to stop protons of specific energies as calculated using SRIM[34]. Calibration was performed using multiple filter configurations to provide several reference points. This involved applying Mylar sheets with thicknesses of 100 μm, 200 μm or 500 μm across the entire MCP, with an example MCP phosphor screen image shown in figure 2(b) (corresponding to the 200 μm case). Each filter thickness imposes a minimum detectable proton energy, $E_{\mathrm{p,min}}$, beyond which protons can reach the detector without being stopped in the Mylar. The values of $E_{\mathrm{p,min}}$, calculated using SRIM simulations[44], were 2.76±0.07 MeV, 4.16±0.10 MeV, and 7.04±0.17 MeV for the respective filters. These values were used as benchmarks to validate the energy calibration and accuracy of the ARISE deflection model.

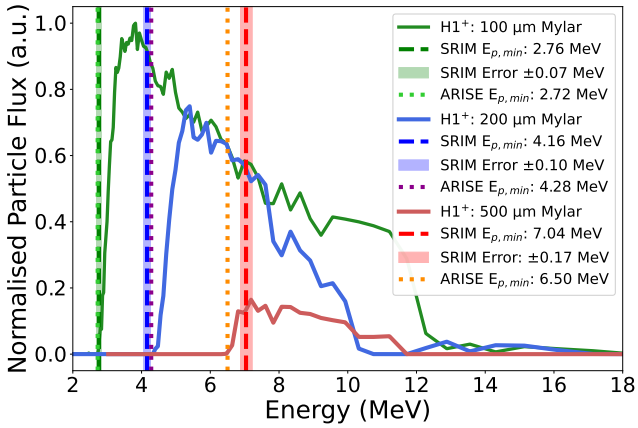

Figure 4 : Measured proton spectra from energy calibration shots. Dashed red, blue and green lines denote the expected minimum transmission energies, $E_{\mathrm{p,min}}$ based on SRIM simulations. Dash-dotted orange, purple and light green lines indicate the corresponding $E_{\mathrm{p,min}}$ values calculated using ARISE. Shaded regions represent uncertainties in the SRIM-derived energy thresholds, defined by the extent of lateral straggling.

As shown in figure 4, there is good agreement between the minimum proton energies $E_{\mathrm{p,min}}$ predicted by ARISE and those expected based on filter thickness. Specifically, ARISE returned values of 2.72 MeV, 4.28 MeV, and 6.50 MeV, compared to the expected values of 2.76±0.07 MeV, 4.16±0.10 MeV, and 7.04±0.17 MeV, respectively. For the thickest filter, there is evidence of increased ion scattering, manifested as a broadening of the apparent minimum energy. Additionally, the nonlinear spacing of energy bins at higher energies, inherent to the TPS geometry, leads to reduced accuracy in the predicted $E_{\mathrm{p,min}}$ in this regime. Nevertheless, across all cases, the predicted values remain in good agreement with those calculated using the SRIM model, confirming the reliability of the ODE-based deflection solver within ARISE.

Following this validation of the modelled energy values, we evaluated several approaches for automatic detection of the maximum proton energy, $E_{\mathrm{p,max}}$. A dataset of 15 randomly selected spectra was extracted from experimentally acquired measurements. For each spectrum, a ground truth value of $E_{\mathrm{p,max}}$ was established through blind manual assessment by four independent researchers, each unaware of the others' selections and the results of any automated method. The ground truth value was defined as the mean of these manual assessments, and the associated standard deviation was used to quantify uncertainty. An automated method is considered well calibrated if its output lies within this confidence interval.

To assess performance, we present the two best-performing $E_{\mathrm{p,max}}$ detection methods and summarise their results against the ground truth data.

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The first of these methods — the less effective of the two — identifies $E_{\mathrm{p,max}}$ by evaluating the local gradient of the ion spectrum and selecting the energy at which the gradient consistently falls below a user-defined threshold. Although conceptually straightforward, this approach is highly sensitive to the chosen threshold parameter, which is somewhat arbitrary and set in the configuration file. As a result, the method is prone to false positives or premature termination, particularly in the presence of noise or weak signal gradients near the high-energy cut-off.

The second method evaluated is based on a least squares regression technique previously applied to time-of-flight ion spectrometers[45]. For a given spectrum, a sliding least squares regression is implemented for successive windows of neighbouring data points along the spectrum. For each window, the gradient and intercept of the best-fit line is computed according to equations (10) and (11):

$$m = \frac{N\sum_{i=1}^{N} x_i y_i - \sum_{i=1}^{N} x_i \sum_{i=1}^{N} y_i}{N\sum_{i=1}^{N} x_i^2 - \left(\sum_{i=1}^{N} x_i\right)^2} \tag{10}$$

$$c = \frac{\sum_{i=1}^{N} x_i^2 \sum_{i=1}^{N} y_i - \sum_{i=1}^{N} x_i \sum_{i=1}^{N} x_i y_i}{N\sum_{i=1}^{N} x_i^2 - \left(\sum_{i=1}^{N} x_i\right)^2} \tag{11}$$

Here, x and y denote the particle energy and particle flux, respectively, and $N = 2r + 1$ is the number of data points included in each local fit, with r being a user-defined parameter specified in the configuration file. The gradient and intercept of each local fit are denoted by m and c, respectively. The maximum proton energy, $E_{\mathrm{p,max}}$, is identified as the point where the final calculated local slope crosses the x-intercept, corresponding to zero flux.

Empirical testing showed that this method performs most reliably for small window sizes, specifically when $r \leq 2$ and $N \leq 5$. This approach offers a key advantage over the fixed-threshold method, as the transition to the background level is determined intrinsically from the spectral shape, rather than through a user-defined parameter.

A comparative analysis of the two methods is presented in figure 5 and table 1.

Figure 5 (a) shows a representative ion spectrum with the ground truth and automatically detected $E_{\mathrm{p,max}}$ values annotated. In this case, the gradient threshold method significantly underperforms relative to the least squares regression, with respect to the ground truth. To explore this further, figure 5(b) presents a comparison of both methods against ground truth values across all 15 analysed spectra. The black dashed line indicates the case of ideal agreement between the ground truth and detected $E_{\mathrm{p,max}}$ values, with the error bars representing the standard deviation in the ground truth value. While neither method is perfectly calibrated, the least squares approach consistently outperforms the threshold method.

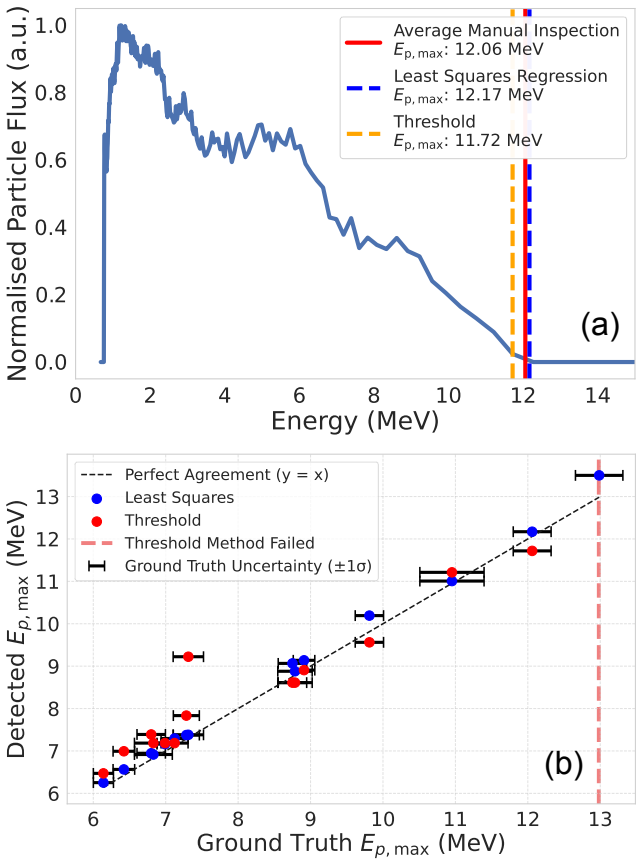

Figure 5 : (a) Example spectrum comparing automatic $E_{\mathrm{p,max}}$ detection using two methods: least squares regression (blue dashed line) and the threshold method (orange dashed line), compared against the ground truth (red solid line). (b) Comparison of $E_{\mathrm{p,max}}$ values across 15 spectra, showing results from least squares regression (blue circles) and the threshold method (red circles) relative to the ground truth. The black dashed line indicates the ideal $y = x$ agreement. Error bars represent standard deviations from the ground truth. A representative case where the threshold method fails is also highlighted (red dashed line).

Table 1 quantitatively compares the two methods using multiple performance metrics, including Mean Absolute Error (MAE), Root Mean Square Error (RMSE), and Mean Absolute Percentage Error (MAPE). Across all metrics, the least squares method outperforms the thresholding approach by more than a factor of two. The 'Within $1\sigma_{\mathrm{human}}$' row further highlights that approximately two-thirds of least squares results fall within one standard deviation of the human-assessed ground truth, compared to just over one-third for the threshold method.

To assess performance in relation to the uncertainty inherent in the ground truth itself, we introduce a normalised metric: the RMSE of each method divided by the mean standard deviation of the human inspection values, $\overline{\sigma}_{\mathrm{human}}$. This dimensionless metric reflects model performance relative to typical human disagreement. A ratio ≤ 1 indicates parity with human reliability; values > 1 suggest inferior accuracy.

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Metric	Least Squares	Threshold
MAE [MeV]	**0.18**	0.41
RMSE [MeV]	**0.22**	0.61
MAPE [%]	**2.06**	5.49
Within $1\sigma_{\text{human}}$	66.7%	35.7%
Failure Rate	0%	6.7%
$\text{RMSE}/\overline{\sigma}_{\text{human}}$	**1.01**	2.82

Table 1 : Performance comparison of $E_{\text{p,max}}$ detection methods. The least squares method significantly outperforms the threshold method, achieving a root mean square error (RMSE) of 0.22 MeV compared to 0.61 MeV. To contextualise these results, we introduce a normalised error metric that compares the automated detection error to the uncertainty in the human-assessed ground truth. Specifically, RMSE values are normalised by the mean standard deviation of the ground truth estimates. Using this metric, the performance of the least squares method is found to be comparable to that of human assessment.

The least squares method yields a value of 1.01, suggesting performance on par with human annotation, while the threshold method returns 2.82, indicating significantly poorer accuracy. Additionally, the gradient threshold method exhibits a higher failure rate and a tendency for 'brittleness' [46], with catastrophic failures observed in certain cases (e.g. no value returned, indicated by the red dashed line in figure 5(b)). This behaviour makes it unsuitable for use in unsupervised data-driven optimisation processes.

Testing ARISE repetition rate - A key aspect of ARISE development is not only the ability to automatically extract ion spectra, but do so at repetition rates exceeding 1 Hz. This capability is essential for real-time optimisation of the ion source, as well as for collecting statistically significant datasets and large-scale training sets for machine learning models.

To validate the maximum effective repetition rate of ARISE, 200 MCP phosphor images from an existing dataset acquired at the SCAPA facility were analysed using ARISE. The analysis was initialised using a prebuilt proton parabola covering the energy range 0.25–20 MeV. The time taken to complete each major processing step in ARISE — background subtraction, multi-threaded median filtering, image handling (including cropping and rotation), automatic ZP detection, and automatic $E_{\text{p,max}}$ extraction — was recorded to identify potential bottlenecks affecting throughput. The computer system used is equipped with an AMD EPYC 9454 processor operating at 2.75 GHz.

When all processing steps are enabled, the average time per image is 292±6 ms, corresponding to a repetition rate of 3.42±0.07 Hz. The dominant time contributions arise from image processing — particularly background subtraction and median filtering — which together account for approximately 83% of the total analysis time. Median filtering was originally implemented to mitigate noise and suppress hot pixels due to hard x-ray hits in earlier datasets. However, more recent data obtained with improved detector shielding indicate that this step is unnecessary. As shown

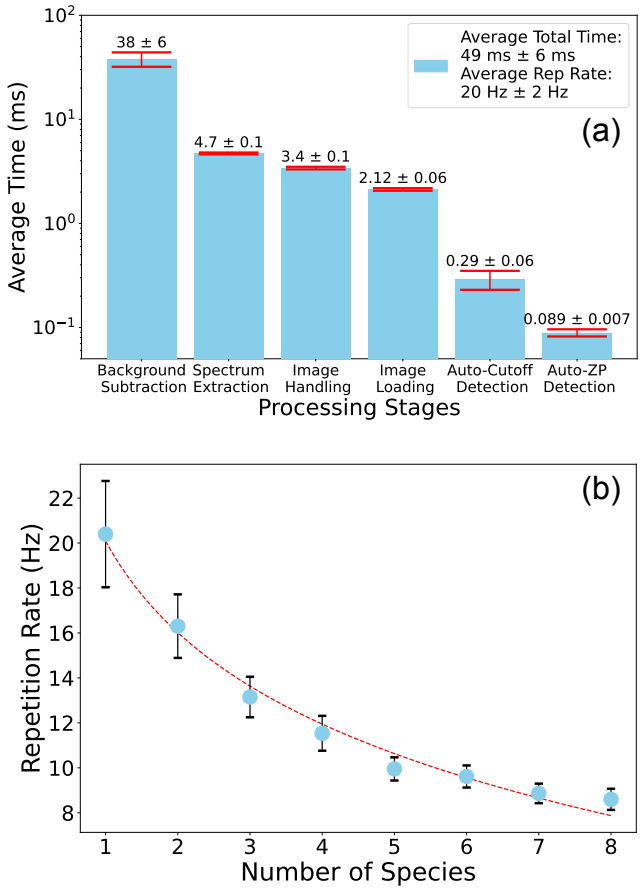

Figure 6 : (a) Bar chart showing the average processing time for key stages of ARISE, with and without median filtering, across 200 data points. The most time-consuming steps are median filtering (when applied) and background subtraction. Without filtering, the pipeline achieves an average repetition rate of 20.40 Hz; applying median filtering reduces this rate significantly to 3.42 Hz. (b) Average repetition rate as a function of the number of ion species, based on 200 shots. Error bars represent one standard deviation in processing time.

in Figure 6(a), when median filtering is omitted, the repetition rate increases significantly to 20±2 Hz, with a mean processing time of 49±6 ms — substantially exceeding the repetition rate of many high-power laser systems and leaving significant allowance for data transfer and the execution of optimisation routines [47].

To evaluate the impact of analysing multiple ion species (e.g., various charge states of carbon and oxygen), the average repetition rate was measured as a function of the number of species processed using the same dataset. Results are shown in figure 6(b). Each species was analysed over the energy range 0.25–20 MeV (total energy, not per nucleon), using 0.1 MeV energy bins. As expected, the repetition rate decreases with increasing species number, falling from its maximum for a single species to 8.60±0.47 Hz when eight species are included. Since image processing is only performed once per image, the overall analysis time scales sub-linearly with the number of species, and the observed

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reduction in repetition rate represents only a factor of ∼2.3.

4.1. Real-time Experimental Demonstration

A real-time experimental validation of ARISE was conducted using the 5 Hz, 350 TW SCAPA laser. Laser pulses of 5.39±0.08 J energy and 29.6±0.4 fs duration were focused to a 1.57±0.04 μm (FWHM) spot onto 10 μm thick steel tape and, in a separate run, onto 13 μm thick Kapton tape targets. During this demonstration the effective repetition rate was limited to 0.2 Hz due to constraints in data transfer speed to the server where data was stored for analysis by ARISE and the readout rate of the diagnostic camera.

In a preliminary demonstration of real-time capability, repeat proton beam measurements were performed without deliberate modification of experimental parameters. Under these stable conditions, ARISE consistently measured $E_{\mathrm{p,max}}$, as shown in figure 7(a), highlighting its ability to actively monitor proton beam properties during periods of stable operation.

To further demonstrate the performance of ARISE in undertaking real-time ion spectra analysis and facilitating data-driven optimisation, the automatically determined $E_{\mathrm{p,max}}$ values were used as the objective function in a Bayesian optimisation feedback loop[48], implemented within our experimental control software[35] and interfaced directly with SCAPA control systems.

The results of the optimisation run are presented in figure 7(b). To initialise the Bayesian optimisation process, five shots were taken with randomised laser energy and pulse duration values, within the ranges 0.28–2.7 J (restricted to mitigate risk of damaging the focusing optic) and 25–135 fs, respectively. The pulse duration was varied using the Group Delay Dispersion (GDD) of a Dazzler[49]. The GDD value corresponding to the shortest pulse duration was 23,000 fs^2. During optimisation, this value was allowed to vary within the range 21,800–24,000 fs^2. Owing to the relationship between GDD and pulse duration, this range enables variation of the pulse duration around its minimum value, and the explored pulse duration range corresponds to a GDD variation of approximately 1,000 fs^2. For each shot, ARISE determined the maximum proton energy, $E_{\mathrm{p,max}}$, which was then used to train an initial Gaussian Process Regression (GPR) model[50], thereby initiating the Bayesian optimisation routine targeting maximisation of $E_{\mathrm{p,max}}$.

In this simple two-parameter optimisation example, figure 7(b) shows that ARISE successfully returns the variation in $E_{\mathrm{p,max}}$ from the random parameter sampling (unfilled symbols) to construct the GPR model that enables the Bayesian optimisation routine to rapidly reach a plateau for maximum laser energy (filled symbols), as expected. Variation in pulse duration variation across the set GDD range was found to have negligible impact on $E_{\mathrm{p,max}}$, consistent with Zimmer *et al.*[51]. This proof-of-principle demonstration of real-time optimisation using only two parameters highlights the capability of ARISE for integration into more complex multi-parameter optimisation frameworks[21,35] and its potential to enable fully automated tuning of laser-driven proton sources in future experimental campaigns.

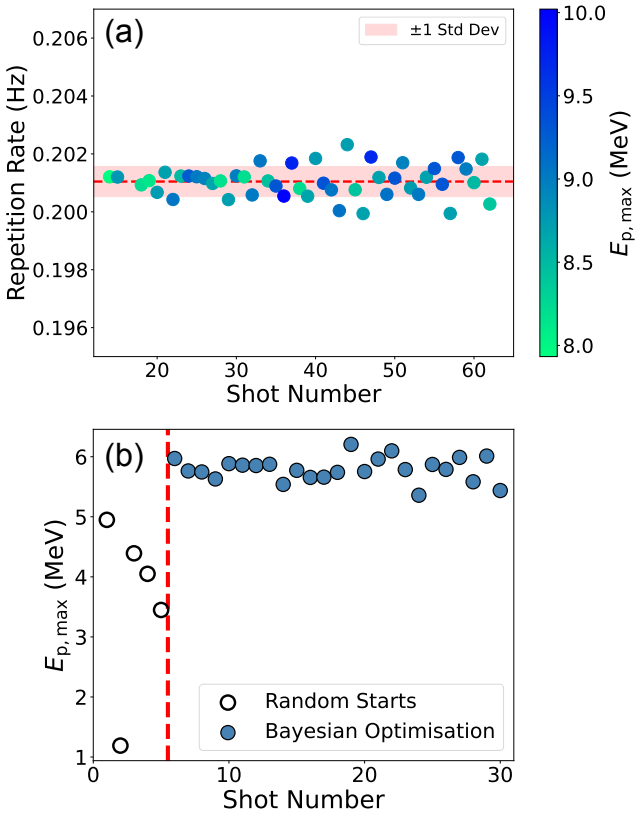

Figure 7 : (a) Example of automatic $E_{\mathrm{p,max}}$ extraction (colour axis) across more than 60 shots at a repetition rate of 0.2 Hz. The shaded region corresponds to 1 standard deviation. (b) Results from an experiment where ARISE-derived $E_{\mathrm{p,max}}$ values were used as the objective function in an open-source Bayesian optimisation feedback loop, in which laser energy and pulse duration were varied by the optimiser. Unfilled symbols represent initial random sampling, while filled symbols correspond to values selected using the Gaussian Process Regression model. The dashed red line separates the random sampling phase from the model-driven optimisation phase.

5. Conclusions

In summary, we have developed and demonstrated ARISE — an algorithm capable of real-time extraction of laser-driven ion spectra at repetition rates exceeding 20 Hz. Its key features include automatic detection of the zero-deflection reference point, background subtraction, and identification of the maximum ion energy. Additionally, it has been deployed to support real-time optimisation of the maximum proton energy via a Bayesian optimisation algorithm. ARISE autonomously extracted $E_{\mathrm{p,max}}$ in real time and this value was used by the Bayesian optimisation algorithm to determine the optimal drive laser parameters. During this experiment, the maximum achievable repetition rate was 0.2 Hz, constrained solely by data transfer speed and the

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diagnostic CCD readout time. Performance testing has demonstrated that ARISE can be applied at multi-Hertz repetition rates.

The development of ARISE represents a significant step towards automated, high-repetition-rate, data-driven optimisation of laser-driven ion sources. Beyond experimental control, it enables the rapid generation of training datasets for neural network–based synthetic diagnostics[26] during live experiments, using the well-established Thomson parabola spectrometer. This capability will facilitate the discovery of new strategies to stabilise and control ion beam properties, advancing progress towards real-world applications[27]. Furthermore, ARISE is easily adaptable to the automated analysis of other charged-particle spectrometers, including electron spectrometers[52] and wide-angle ion spectrometers[53], promoting broader adoption of data-driven approaches in laser–plasma diagnostics.

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