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Contents

Co	ontents	V
Ι	A FLEXIBLE PYTHON TOOL FOR FOURIER-TRANSFORM NOISE SPECTROSCOPY	1
1	Introduction	3
2	Theory of spectral noise estimation 2.1 Spectrum estimation from time series	5 5 7 8
3	The python_spectrometer software package 3.1 Package design and implementation	11 11 13 14 14 17
4	Conclusion and outlook	19
II	CHARACTERIZATION AND IMPROVEMENTS OF A MILLIKELVIN CONFOCAL MICROSCOPE	21
III	ELECTROSTATIC TRAPPING OF EXCITONS IN SEMICONDUCTOR MEMBRANES	23
IV	A FILTER-FUNCTION FORMALISM FOR UNITAL QUANTUM OPERATIONS	25
Aı	PPENDIX	27
Bil	bliography	29
Lis	st of Terms	31

Part I

A FLEXIBLE PYTHON TOOL FOR FOURIER-TRANSFORM NOISE SPECTROSCOPY

oise is ubiquitous in condensed matter physics experiments, and in mesoscopic systems in particular it can easily drown out the sought-after signal. Hence, characterizing (and subsequently mitigating) noise is an essential task for the experimentalist. But noise comes in as many different forms as there are types of signal sources and detectors, whether it be a voltage source or a photodetector, and while some instruments have built-in solutions for noise analysis, they vary in functionality and capability. Moreover, the measured signal often does not directly correspond to the noisy physical quantity of interest, making it desirable to be able to manipulate the raw data before processing.

Theory of spectral noise estimation



HERE exists a multitude of methods for estimating noise properties of a classical signal x(t).

If the noisy process x(t) has Gaussian statistics, meaning that the value at a given point in time follows a normal distribution with some mean μ and variance σ^2 over multiple realizations of the process, it can be fully described by the power spectral density (PSD) $S(\omega)$. For the purpose of noise estimation, the assumption of Gaussianity is a rather weak one as the noise typically arises from a large ensemble of individual fluctuators and is therefore well approximated by a Gaussian distribution by the central limit theorem. Even if the process x(t) is not perfectly Gaussian, non-Gaussian contributions can be seen as higher-order contributions if viewed from the perspective of perturbation theory, and therefore the PSD still captures a significant part of the statistical properties. For this reason, the PSD is the central quantity of interest in noise spectroscopy and I will discuss some of its properties in the following.

For real signals $x(t) \in \mathbb{R}$, $S(\omega)$ is an even function and one therefore distinguishes the *two-sided* PSD $S^{(2)}(\omega)$ defined over \mathbb{R} from the *one-sided* PSD $S^{(1)}(\omega) = 2S^{(2)}(\omega)$ defined only over \mathbb{R}^+ . Complex signals $x(t) \in \mathbb{C}$ such as those generated by Lock-in amplifiers after demodulation in turn have asymmetric, two-sided PSDs.

2.1 Spectrum estimation from time series

To see how the PSD may be estimated from time-series data, consider a continuous wide-sense stationary⁴ signal in the time domain $x(t) \in \mathbb{C}$ that is observed for some time T. We define the windowed Fourier transform of x(t) and its inverse by⁵

$$\hat{x}_T(\omega) = \int_0^T dt \, x(t) e^{-i\omega t} \tag{2.1}$$

and
$$x(t) = \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \hat{x}_T(\omega) e^{i\omega t},$$
 (2.2)

i.e., we assume that outside of the window of observation x(t) is zero. The auto-correlation function of x(t) is given by

$$C(\tau) = \langle x(t)^* x(t+\tau) \rangle \tag{2.3}$$

$$= \lim_{T \to \infty} \frac{1}{T} \int_{0}^{T} dt \, x(t)^{*} x(t+\tau), \tag{2.4}$$

where $\langle \cdot \rangle$ is the ensemble average over multiple realizations of the process and the last equality holds true for ergodic processes. Expressing x(t) in terms of its Fourier representation (Equation 2.1) and reordering the integrals, we get⁶

1: We discuss only classical noise here, meaning x(t) commutes with itself at all times. For descriptions of and spectroscopy protocols for quantum noise refer to Refences 1 and 2, for example.

lay out some others

2: The term *power spectrum* is often used interchangably. I will do so as well, but emphasize at this point that in digital signal processing in particular, the *spectrum* is a different quantity from the *spectral density*. See also sidenote 17 in Chapter 3.

maybe a classical signal processing ref?

3: As an example, consider electronic devices, where voltage noise arises from a large number of defects and other charge traps in oxides being populated and depopulated at certain rates γ . The ensemble average over these so-called two-level fluctuators (TLFs) then yields the well-known 1/f-like noise spectra (at least for a large density [3]).

flesh out this sidenote?

flesh out, organize

4: For a wide-sense stationary (also called weakly stationary) process x(t), the mean is constant and the auto-correlation function $C(t,t') = \langle x(t)^*x(t') \rangle$ is given by $\langle x(t)^*x(t+\tau) \rangle = \langle x(0)^*x(\tau) \rangle$ with $\tau = t' - t$. That is, it is a function of only the time lag τ and not the absolute point in time. For Gaussian processes as discussed here, this also implies stationarity [4]. The property further implies that $C(\tau)$ is an even function.

sketch of auto-correlation function?

5: In this chapter we will always denote the Fourier transform of some quantity ξ using the same symbol with a hat, $\hat{\xi}$.

6: Mathematicians might at this point argue the integrability of x(t), but as we deal with physical processes with finite bandwidth (and have no shame), we do not

$$C(\tau) = \lim_{T \to \infty} \frac{1}{T} \int_0^T dt \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \hat{x}_T(\omega)^* e^{-i\omega t} \int_{-\infty}^{\infty} \frac{d\omega'}{2\pi} \hat{x}_T(\omega') e^{i\omega'(t+\tau)}$$
(2.5)

$$= \lim_{T \to \infty} \frac{1}{T} \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \int_{-\infty}^{\infty} \frac{d\omega'}{2\pi} \hat{x}_{T}(\omega)^{*} \hat{x}_{T}(\omega') e^{i\omega'\tau} \int_{0}^{T} dt \, e^{it(\omega'-\omega)}$$
 (2.6)

7: Note that, because x(t) is wide-sense stationary, we may shift the limits of integration $\int_0^T \to \int_{-T/2}^{+T/2}$.

The innermost integral approaches a δ -function for large T, allowing us to further simplify this under the limit as

$$C(\tau) = \lim_{T \to \infty} \frac{1}{T} \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \int_{-\infty}^{\infty} \frac{d\omega'}{2\pi} \hat{x}_{T}(\omega)^{*} \hat{x}_{T}(\omega') e^{i\omega'\tau} \delta(\omega' - \omega)$$
 (2.7)

$$= \lim_{T \to \infty} \frac{1}{T} \int_{-\infty}^{\infty} \frac{\mathrm{d}\omega}{2\pi} |\hat{x}_T(\omega)|^2 \mathrm{e}^{\mathrm{i}\omega\tau}$$
 (2.8)

$$= \int_{-\infty}^{\infty} \frac{\mathrm{d}\omega}{2\pi} S(\omega) \mathrm{e}^{\mathrm{i}\omega\tau} \tag{2.9}$$

with the PSD

$$S(\omega) = \lim_{T \to \infty} \frac{1}{T} |\hat{x}_T(\omega)|^2 \tag{2.10}$$

$$= \int_{-\infty}^{\infty} d\tau C(\tau) e^{-i\omega\tau}$$
 (2.11)

Equation 2.9 is the Wiener-Khinchin theorem that states that the auto-correlation function $C(\tau)$ and the PSD $S(\omega)$ are Fourier-transform pairs [4]. Furthermore, defining the latter through Equation 2.10 gives us an intuitive picture of the PSD if we recall Parseval's theorem,

$$\int_{-\infty}^{\infty} \frac{\mathrm{d}\omega}{2\pi} \frac{1}{T} |\hat{x}_T(\omega)|^2 = \frac{1}{T} \int_{-\infty}^{\infty} \mathrm{d}t |x(t)|^2. \tag{2.12}$$

That is, the total power P contained in the signal x(t) is given by integrating over the PSD. Similarly, the power contained in a band of frequencies $[\omega_1, \omega_2]$ is given by

$$P(\omega_1, \omega_2) = \text{RMS}_S(\omega_1, \omega_2)^2 \tag{2.13}$$

$$= \int_{\omega_1}^{\omega_2} \frac{\mathrm{d}\omega}{2\pi} S(\omega) \tag{2.14}$$

where $\mathrm{RMS}_S(\omega_1,\omega_2)$ is the root mean square within this frequency band. These relations are helpful when analyzing noise PSDs to gauge the relative weight of contributions from different frequency bands to the total noise power.

Equation 2.10 represents the starting point for the experimental spectrum estimation procedure. Instead of a continuous signal $x(t), t \in [0, T]$, so we only discuss the problem of equally spaced samples here. Variants for

$$x_n, \quad n \in \{0, 1, \dots, N-1\}$$
 (2.15)

defined at times $t_n = n\Delta t$ with $T = N\Delta t$ and where $\Delta t = f_{\rm s}^{-1}$ is the sampling interval (the inverse of the sampling frequency $f_{\rm s}$). Invoking the ergodic theorem, we can replace the long-term average in Equation 2.10 by the ensemble average over M realizations i of the noisy signal $x_n^{(m)}$ and

8: We only discuss the problem of equally spaced samples here. Variants for spectral estimation of time series with unequal spacing exist [5, 6].

write

$$S_n = \frac{1}{M} \sum_{i=0}^{M-1} \left| \hat{x}_n^{(m)} \right|^2 \tag{2.16}$$

$$=\frac{1}{M}\sum_{i=0}^{M-1}S_n^{(m)}\tag{2.17}$$

where $\hat{x}_n^{(m)}$ is the discrete Fourier transform of $x_n^{(m)}$, we defined the *periodogram* of $x_n^{(m)}$ by

$$S_n^{(m)} = \left|\hat{x}_n^{(m)}\right|^2,$$
 (2.18)

and S_n is an *estimate* of the true PSD sampled at the discrete frequencies $\omega_n=2\pi n/T\in 2\pi\times\{-f_{\rm s}/2\,,\ldots,\,f_{\rm s}/2\}.^9$ Equation 2.16 is known as Bartlett's method [7] for spectrum estimation. ¹⁰

To better understand the properties of this estimate, let us take a look at the parameters Δt , N, and M. The sampling interval Δt defines the largest resolvable frequency by the Nyquist sampling theorem,

$$f_{\text{max}} = \frac{f_{\text{s}}}{2} = \frac{1}{2\Delta t}.$$
 (2.19)

In turn, the number of samples N determines the frequency resolution Δf , or smallest resolvable frequency,

$$f_{\min} = \Delta f = \frac{1}{T} = \frac{1}{N\Delta t} = \frac{f_s}{N}.$$
 (2.20)

Lastly, M determines the variance of the set of periodograms $\left\{S_n^{(m)}\right\}_{i=0}^{M-1}$ and hence the accuracy of the estimate S_n .

In practice, the ensemble realizations i are of course obtained sequentially, implying that one acquires a time series of data $x_n, n \in \{0, 1, ..., NM-1\}$ and partitions these data into M sequences of length N. It becomes clear, then, that the Bartlett average (Equation 2.16) trades spectral resolution (larger N) for estimation accuracy (larger M) given the finite acquisition time $T = NM\Delta t$.

An improvement in data efficiency can be obtained using Welch's method [8]. To see how, we first need to discuss spectral windowing.

2.2 Window functions

Partitioning a signal x_n into M sections $x_n^{(m)}$ of length N is mathematically equivalent to multiplying the signal with the rectangular window function given by 11

$$w_n^{(m)} = \begin{cases} 1 & \text{if } (m-1)N \le n < mN \text{ and} \\ 0 & \text{else} \end{cases}$$
 (2.21)

so that $x_n^{(m)}=x_nw_n^{(m)}$. Now recall that multiplication and convolution are duals under the Fourier transform, implying that

$$\hat{x}_n^{(m)} = \hat{x}_n * \hat{w}_n^{(m)}, \tag{2.22}$$

9: We blithely disregard integer algebra issues occuring here for conciseness and leave it as an exercise for the reader to figure out what the exact bounds of the set of ω_n are.

10: By taking the limit $M \to \infty$ one recovers the true PSD,

$$\lim_{M\to\infty} S_n = S(\omega_n).$$

The continuum limit is as always obtained by sending $\Delta t \rightarrow 0, N \rightarrow \infty, N\Delta t = \text{const.}$

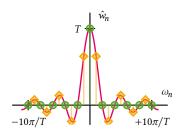


Figure 2.1: The Fourier representation of the rectangular window in continuous time (solid line) and for discrete frequencies $\omega_n = 2\pi n/T$ (circles). Introducing a phase shift, that is, shifting the window with respect to the signal in time, effectively shifts $\omega_n \to \omega_{n+\eta}$ as indicated for $\eta = 1/2$ (diamonds). This incurs scalloping loss.

11: This window is also known as the boxcar or Dirichlet window.

12: $\operatorname{sinc}(x) = \sin(x)/x$.

where the Fourier representation of the rectangular window 12

$$\hat{\mathbf{w}}_n^{(m)} = \hat{\mathbf{w}}_n e^{-i(m-1/2)\omega_n T}, \tag{2.23}$$

$$\hat{w}_n = T \operatorname{sinc}\left(\frac{\omega_n T}{2}\right). \tag{2.24}$$

Figure 2.1 shows the unshifted rectangular window \hat{w}_n in Fourier space. We can hence understand the Fourier spectrum of $x_n^{(m)}$ as sampling \hat{x}_n with the probe $\hat{w}_n^{(m)}$. However, while in the continuum limit (c.f. sidenote 10) Equation 2.24 tends towards $\delta(\omega_n)$ and thus will produce a faithful reconstruction of the true spectrum, the finite frequency spacing Δf of discrete signals and finite observation length T introduce a finite bandwidth of the probe as well as sidelobes. These effects induce what is known as spectral leakage and scalloping loss [4, 9] and lead to artifacts and deviations of the spectrum estimator S_n from the true spectrum $S(\omega_n)$.

For this reason, a plethora of *window functions* have been introduced to mitigate the effects of spectral leakage. Key properties of a window are the spectral bandwidth (center lobe width) and sidelobe amplitude between which there typically is a tradeoff.¹³

A window frequently used in spectral analysis is the Hann window [11],

$$w_n^{(m)} = \begin{cases} \cos^2\left(\frac{\pi n}{N}\right) & \text{if } (m-1)N \le n < mN \text{ and} \\ 0 & \text{else} \end{cases}$$
 (2.25)

with the Fourier representation of the unshifted window,

$$\hat{w}_n = T \operatorname{sinc}\left(\frac{\omega_n T}{2}\right) \times \frac{1}{2(1 - \omega_n T/2\pi)(1 + \omega_n T/2\pi)},\tag{2.26}$$

shown in Figure 2.2. The favorable properties of the Hann window are apparent when compared to the rectangular window in Equation 2.24 and shown in Figure 2.1; the sidelobes are quadratically suppressed while the center lobe is only slightly broadened.

Another favorable property of the Hann window is that $w_0^{(0)} = w_{N-1}^{(0)} = 0$. This suppresses detrimental effects arising from a possible discontinuity $(x_0^{(0)} \neq x_{N-1}^{(0)})$ at the edge of a data segment related to the discrete Fourier transform, which assumes periodic data.¹⁴

2.3 Welch's method

Contemplating Equation 2.25, one might come to the conclusion that using a window such as this is not very data efficient in the sense that a large fraction of samples located at the edge of the window is strongly suppressed and hence does not contribute significantly to the spectrum estimate. To alleviate this lack of efficiency, one can introduce an overlap between adjacent data windows. That is, instead of partitioning the data x_n into M non-overlapping sections of length N, one shifts the mth window forward by -mK with K>0 the overlap. Finally, the periodogram (Equation 2.18) is computed for each window and subsequently averaged to obtain the spectrum estimator (Equation 2.16).

This method of spectrum estimation is known as Welch's method [8]. One can show [8] that the correlation between the periodograms of adjacent, overlapping windows is sufficiently small to avoid a biased estimate. The

13: Wikipedia gives a good overview of existing window functions [10].

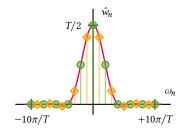
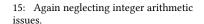


Figure 2.2: The Fourier representation of the Hann window in continuous time (solid line) and for discrete frequencies ω_n (circles). Diamonds indicate discrete sampling when the window completely out of phase with the signal (c. f. Figure 2.1).

14: Although this can usually also be achieved approximately by detrending the data before performing the Fourier transform, which is a good idea in any case.

overlap naturally depends on the choice of window; a typical value for the Hann window K=N/2 with which one would obtain M=2L/N-1 windows for data of length L.¹⁵



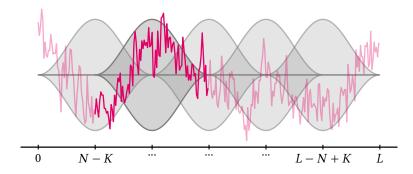


Figure 2.3: Illustration of Welch's method for spectrum estimation. The data (pink) of length L is partitioned into K=2L/N-1 segments of length N. Each segment is multiplied with a window function (gray) which reduces spectral leakage and other artifacts. A finite overlap K between adjacent windows (gray) ensures efficient sample use.

Figure 2.3 conceptually illustrates Welch's method for a trace of 1/f noise with L=300 samples in total. Choosing the Hann window and an overlap of 50 % results in M=5 segments for a window length of N=100. The data in the second window is highlighted.

2.3.1 Parameters

We are now in a position to discuss how the various parameters of a time series relate to both to physical parameters of the resulting spectrum estimate and to each other. To this end, we will go through the typical procedure of acquiring a spectrum estimate using Welch's method chronologically.

To acquire data using some form of (digital) data acquisition device (DAQ), one usually needs to specify two parameters first: the total number of samples to be acquired, L, and the sample rate, f_s . This results in a measurement of duration $T = L\Delta t$ where $\Delta t = f_s^{-1}$ as previously mentioned. The choice of f_s already induces an upper bound on the first parameter characterizing the PSD estimate: the largest resolvable frequency $f_{\rm max} \leq f_s/2$ (c.f. Equation 2.19, but note that we allow $f_{\rm max}$ to be smaller than half the sample rate in anticipation of hardware constraints). Next, we choose a number of Welch averages, M, i.e., data partitions, and their overlap, K. In doing so, one fixes the number of samples per partition N and thereby induces the lower bound on the second parameter characterizing the PSD estimate: the frequency spacing $\Delta f = 1/N \leq f_{\rm min}$ (c.f. Equation 2.20). The sample of the parameter of the property of the property of the property of the property of the parameter of the property of the property of the parameter of the parameter

Table 2.1: Overview of spectrum estimation parameters. The parameters can be assigned into three groups: 1. DAQ parameters configuring the data acquisition device, 2. Welch parameters specifying the periodogram averaging, and 3. Spectrum properties induced by the above.

1. DAQ parameters

L Total number of samples

 $f_{\rm s}$ Sample rate

2. Welch parameters

K Number of overlap samples

N Number of segment samples

M Number of Welch segments

3. Spectrum parameters

 $\begin{array}{ll} f_{\min} & \text{Smallest resolvable frequency} \\ f_{\max} & \text{Largest resolvable frequency} \end{array}$

16: Technically, the smallest resolvable frequency in a fast Fourier transform (FFT) is zero, of course. But as data is typically detrended (a constant or linear trend subtracted) before computation of the periodogram, the smallest *meaning-ful* frequency is given by f_{\min} .

D(A)G for parameter interdependencies?

The python_spectrometer software package



N to

N this chapter, I will lay out the design and functionality of the ${\tt python_spectrometer}$ Python package. 1

3.1 Package design and implementation

The python_spectrometer package provides a central class, Spectrometer, that users interact with to perform data acquisition, spectrum estimation, and plotting. It is instantiated with an instance of a child class of the DAQ base class that implements an interface to various DAQ hardware devices. New spectra are obtained by calling the Spectrometer.take() method with all acquisition and metadata settings.

In the following, I will go over the design of these aspects of the package in more detail.

3.1.1 Data acquisition

The daq module contains on the one hand the declaration of the DAQ abstract base class and its child class implementations, and on the other the settings module, which defines the DAQSettings class. This class is used in the background to validate data acquisition settings both for consistency (c.f. Subsection 2.3.1) and hardware constraints.

To better understand the necessity of this functionality, consider the typical scenario of a physicist² in the lab. Alice has wired up her experiment, performed a first measurement, and to her dismay discovered that the data is too noisy to see the sought-after effect. She sets up the python_1 spectrometer code to investigate the noise spectrum of her measurement setup. From her noisy data she could already estimate the frequency of the most harrowing noise, so she knows the frequency band $[f_{\min}, f_{\max}]$ she is most interested in. But because she is lazy,³ she does not want to do the mental gymnastics to convert f_{\min} to the parameter that her DAQ device understands, L (see Table 3.1), especially considering that L depends on the number of Welch averages and the overlap. Furthermore, while she could just about do the conversion from $f_{
m max}$ to the other relevant DAQ parameter, f_s , in her head, her device imposes hardware constraints on the allowed sample rates she can select! The DAQSettings class addresses these issues. It is instantiated with any subset of the parameters listed in Table 3.14 and attempts to resolve the parameter interdependencies $\stackrel{-}{\text{lined out in Subsection 2.3.1 upon calling DAQSettings.to_consistent}}_{-1}$ dict().⁵ This either infers those parameters that were not given from those that were or, if not possible, uses a default value. Child classes of the DAQ class can subclass DAQSettings to implement hardware constraints such as a finite set of allowed sampling rates or a maximum number of samples per data buffer.

For instance, Alice might want to measure the noise spectrum in the frequency band [1.5 Hz, 72 kHz]. Although she would not have to do this explicitly,⁶ she could inspect the parameters after resolution using the code shown in Listing 3.1.

1: The package repository is hosted on GitLab. Its documentation is automatically generated and hosted on GitLab as well. Releases are automatically published to PyPI and allow the package to be installed using pip install python-spectrometer.

Table 3.1: Variable names used in Chapter 2 and their corresponding parameter names as used in python_spectrometer and scipy.signal.welch() [12].

Variable	Parameter
L	n_pts
$f_{ m s}$	fs
K	noverlap
N	nperseg
M	n_seg
f_{\min}	f_min
$f_{ m max}$	f_max

- 2: Let's call her Alice.
- 3: Physicists generally are.

- 4: DAQSettings inherits from the builtin dict and as such can contain arbitrary other keys besides those listed in Table 3.1. However, automatic validation of parameter consistency is only performed for these special keys.
- 5: Since the graph spanned by the parameters is not acyclic, this only works *most* of the time.
- 6: Settings are automatically parsed when passed to the take() method of the Spectrometer class.

Listing 3.1: DAQSettings example show-casing automatic parameter resolution. n_avg determines the number of outer averages, *i. e.*, the number of data buffers acquired and processed individually.

```
>>> from python_spectrometer.daq import DAQSettings
>>> settings = DAQSettings(f_min=1.5, f_max=7.2e4)
>>> settings.to_consistent_dict()
{'f_min': 1.5,
   'f_max': 72000.0,
   'fs': 144000.0,
   'df': 1.5,
   'nperseg': 96000,
   'noverlap': 48000,
   'n_seg': 5,
   'n_pts': 288000,
   'n_avg': 1}
```

```
{'f_min': 14.30511474609375,
  'f_max': 72000.0,
  'fs': 234375.0,
  'df': 14.30511474609375,
  'nperseg': 16384,
  'noverlap': 0,
  'n_seg': 1,
  'n_pts': 16384,
  'n_avg': 1}
```

Listing 3.2: Resolved settings for the same input parameters as in Listing 3.1 but for the ZurichInstrumentsMFLIScope backend with hardware constraints on n_pts and fs

[13]: (n.d.), Scope Module - LabOne API User Manual

7: And issued a warning to inform the user their requested settings could not be matched.

Listing 3.3: DAQ workflow pseudocode. A MyDAQ object (representing the instrument My) is instantiated with a driver object (for instance a QCoDeS Instrument). The instrument is configured with the given user_settings. Calling the generator function daq.acquire() with the actual device settings returns a generator, iterating over which yields one data buffer per iteration. The data buffers can then be passed to further processing functions (the PSD estimator in our example).

- 8: Which might differ from the requested settings as outlined above.
- 9: $\it I.e.$, the number of time series data batches acquired, as opposed to the number of Welch averages n_seg within one batch.
- 10: C.f. the implementation of the AlazarTech ATS9440 digitizer card.

If the instrument she'd chosen for data acquisition had been a Zurich Instruments MFLI's "Scope" module [13], the same requested settings would have resolved to those shown in Listing 3.2.⁷ This is because the Scope module constrains $L \in [2^{12}, 2^{14}]$ and $f_{\rm s} \in 60\,{\rm MHz} \times 2^{[-16,0]} \approx \{915.5\,{\rm Hz}, \dots, 30\,{\rm MHz}, 60\,{\rm MHz}\}.$

As already mentioned, the DAQ base class implements a common interface for different hardware backends, allowing the Spectrometer class to be hardware agnostic. That is, changing the instrument that is used to acquire the data does not necessitate adapting the code used to interact with the Spectrometer. To enable this, different instruments require small wrapper drivers that map the functionality of their actual driver onto the interface dictated by the DAQ class. This is achieved by subclassing DAQ and implementing the DAQ.setup() and DAQ.acquire() methods. Their functionality is best illustrated by the internal workflow as representatively shown in Listing 3.3.

```
daq = MyDAQ(driver_handle)

parsed_settings = daq.setup(**user_settings)
acquisition_generator = daq.acquire(**parsed_settings)

for data_buffer in acquisition_generator:
    estimate_psd(data_buffer)
```

When acquiring a new spectrum, all settings supplied by the user are first fed into the setup() method where instrument configuration takes place. The method returns the actual device settings, which are then forwarded to the acquire() generator function. Here, the instrument is armed (if necessary), and subsequently data is fetched from the device and yielded to the caller n_avg times, where n_avg is the number of outer averages. An exemplary implementation of a DAQ subclass for a fictitious instrument is shown in Listing 3.4. In addition to the methods to configure the instrument and perform data acquisition, it is possible to override the DAQSettings property to implement instrument-specific hardware constraints such as, in this example, the number of samples per buffer being constrained to the discrete interval [1, 2048]. Leveraging the qutil.domains module, more complex constraints such as sample rates restricted to an internal clock rate divided by a power of two can be specified.

```
# daq/mydaq.py
import dataclasses
from qutil.domains import DiscreteInterval
from .base import DAQ
from .settings import DAQSettings
@dataclasses.dataclass
class MyDAQ(DAQ):
    handle: mydriver.DeviceHandle
    @property
    def DAQSettings(self) -> type[DAQSettings]:
       class MyDAQSettings(DAQSettings):
           ALLOWED_NPERSEG = DiscreteInterval(1, 2048)
        return MyDAQSettings
    def setup(self, **settings) -> dict:
        settings = self.DAQSettings(settings)
        parsed_settings = settings.to_consistent_dict()
        self.handle.configure(parsed_settings)
        return parsed_settings
    def acquire(self, n_avg: int, *, **settings) -> Generator:
       self.handle.arm(n_avg)
        for _ in range(n_avg):
            self.handle.wait_for_trigger()
           yield self.handle.fetch()
        return self.handle.metadata
```

Listing 3.4: Exemplary code for a DAQ implementation of some instrument with given driver class DeviceHandle in the package mydriver. The MyDAQ class is instantiated with a DeviceHandle instance. Optionally, the DAQSettings property can be overridden to implement hardware constraints or default values for data acquisition parameters. For this, the gutil.domains module provides several classes that represent bounded domains and sets. The setup() method parses the given acquisition settings and configures the instrument through the external driver interface handle. configure(). The acquire() method arms the instrument (if necessary) and loops over the number of outer averages, n_avg. In the body of the loop, it can wait for external triggers (or send software triggers) before yielding a batch of data fetched from the external driver interface. Once acquisition is done, the method can return arbitrary metadata to the Spectrometer object to attach to the stored data.

3.1.2 Data processing

Once time series data has been acquired using a given DAQ backend, it could in principle immediately be used to estimate the PSD following Equation 2.16. However, it is often desirable to transform, or process, the data in some fashion. This can include simple transformations such as accounting for the gain of a transimpedance amplifier (TIA) and convert the voltage back to a current, ¹¹ or more complex ones such as applying calibrations. In particular, since the process of computing the PSD already involves Fourier transformation, the processing can also be performed in frequency space.

In python_spectrometer, this can be done using a procfn (in the time domain) or fourier_procfn (in the Fourier domain). The former is specified as an argument directly to the Spectrometer constructor. It is a callable with signature (x, **kwargs) -> xp, that is, takes the time series data as its first (positional) argument and arbitrary settings that are passed through from the take() method as keyword arguments, and returns the processed data. Listing 3.5 shows a simple function that accounts for the gain of an amplifier.

The latter is specified in the psd_estimator argument of the Spectrometer constructor. This argument allows the user to specify a custom estimator for the PSD, in which case a callable is expected. Otherwise, it should be a mapping containing parameters for the default PSD estimator, <code>scipy.jsignal.welch()</code> [12]. Here, the keyword fourier_procfn should be a callable with signature (xf, f, **kwargs) -> (xfp, fp). That is, it should take the frequency-space data, the corresponding frequencies, and arbitrary keyword arguments and return a tuple of the processed data and the corresponding frequencies. The latter are required in case the

Introduce Bob?

11: Although it is of course less than trivial to discriminate between current and voltage noise in a TIA.

```
def comp_gain(x, gain=1.0, **_):
    return x / gain
```

Listing 3.5: A simple proofn, which converts amplified data back to the level before amplification. More complex processing chains can concisely be defined with qutil.functools.FunctionChain that pipes the output of one function into the input of the next.

[12]: (n.d.), Welch - SciPy v1.15.2 Manual
12: I.e., the psd_estimator argument would be {"fourier_procfn": fn}.

13: One example is the octave_ $_{J}$ band_rms() function from the qutil. $_{J}$ signal_processing module [14]

```
def derivative(xf, f, n=0, **_):
    return xf / (2j * pi * f)**n
```

Listing 3.6: A simple fourier_procfn, which calculates the (anti-)derivative.

ref

function modifies the frequencies. ¹³ A simple example for a processing function in Fourier space is shown in Listing 3.6, which computes the (anti-)derivative of the data using the fact that

$$\frac{\partial^n}{\partial t^n} \xrightarrow{\text{F.T.}} (i\omega)^n \tag{3.1}$$

under the Fourier transform. In Part II, I discuss more complex use-cases of the processing functionality included in python_spectrometer in the context of vibration spectroscopy.

3.2 Feature overview

überleitung

3.2.1 Serial spectrum acquisition

Now that we have a basic understanding of the design choices underlying python_spectrometer, let us discuss the typical workflow of using the package. The default mode for spectrum acquisition using python_j spectrometer revolves around the take() method. Key to this workflow is the idea that each acquired spectrum can be assigned a comment that allows to easily identify a spectrum in the main plot. For instance, this comment could contain information about the particular settings that were active when the spectrum was recorded, or where a particular cable was placed.

Consider as an example the procedure of "noise hunting", *i.e.*, debugging a noisy experimental setup. The experimentalist, 14 having discovered that his data is noisier than expected, sets up the Spectrometer class with an instance of the DAQ subclass for the DAQ instrument connected to his sample, a Zurich Instruments MFLI. 15 Choosing the frequency bounds, say $f_{\rm min}=10\,{\rm Hz}$ and $f_{\rm max}=100\,{\rm kHz}$, and using the sensible defaults for the remaining spectrum parameters, Bob first grounds the input of his DAQ to record a *baseline* spectrum. Thus far, his code would hence look something like that shown in Listing 3.7, which produces the plot shown in Figure 3.1.

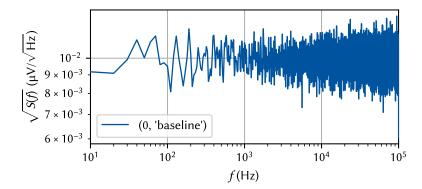
After acquiring the baseline, he next ungrounds the DAQ to obtain a representative spectrum of the noise in an actual measurement. He then proceeds by tweaking things on his setup, testing out different parameters, *etc.* Every time he changes something, he acquires another spectrum using take(), labeling each with a meaningful comment for identification.

This leaves him with the spectrometer plot as shown in Figure 3.2. While working, Bob realizes he'd like see the signal in the time-domain as well. He easily achieves this by setting spect.plot_timetrace = True,

14: Let's call him Bob.

15: For the MFLI, DAQ subclasses for both the Scope and the DAQ module are implemented. The former works with data before and the latter with data after demodulation.

Listing 3.7: Setup and serial workflow using the python_spectrometer package. session and device are Application Programming Interface (API) objects of the zhinst.toolkit driver package. It is therefore possible to simply use the driver objects that are already in use in the measurement setup. The procfn and processed_unit arguments help converting raw data into a more human-friendly unit.



```
settings = {'f_min': 1e1, 'f_max': 1e5, 'daq': 'connected'}
spect.take('connected', **settings)
spect.take('lifted cable', cable='lifted', **settings)
spect.take('jumped', **settings)
```

which adds an oscilloscope subplot to spectrometer figure as shown in Figure 3.3.

Bob now observes that the noise spectra he has recorded display many sharp peaks in particular at high frequencies while the 1/f noise floor seems pretty consistent across different measurements. This makes it harder for him to evaluate whether any of his changes are actually an improvement or not. The python_spectrometer package allows addressing this by plotting the integrated spectra in another subplot. Bob's spectrometer figure after setting spect.plot_cumulative = True is shown in Figure 3.4. In the case that spect.plot_amplitude == True, this new subplot shows the root mean square (RMS) in the band $[f_{\min}, f]$,

$$RMS_{S}(f) \equiv RMS_{S}(f_{min}, f), \tag{3.2}$$

and the band power (Equation 2.13) otherwise.

The cumulative RMS plot already helps, but Bob would like a more quantitative comparison of relative spectral powers. Therefore, he rescales the spectra in terms of their relative powers expressed in dB¹⁶ by applying the following settings, which produces the plot shown in Figure 3.5:

```
spect.plot_dB_scale = True
spect.plot_amplitude = False
spect.plot_density = False
```

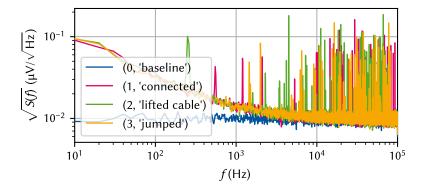


Figure 3.1: The python_spectrometer plot after acquiring the (here: synthetic) baseline spectrum. By default, the amplitude spectral density (ASD) = √PSD is displayed in the main plot. Each spectrum is assigned a unique identifier key consisting of an incrementing integer and the user comment, and can be refered to by either (or both) when interacting with the object.

Listing 3.8: Code to acquire additional spectra. Arbitrary key-value pairs can be passed to the take() method, which are stored as metadata if they do not apply to any functions downstream in the data processing chain.

16: Recall that the decibel is defined by the ratio L_P of two powers P_1 , P_2 as [15]

$$L_P = 10 \log_{10} \left(\frac{P_1}{P_2}\right) \, \mathrm{dB}.$$

Figure 3.2: The python_spectrometer plot after acquiring additional (synthetic) spectra. Each spectrum is uniquely identified by a two-tuple of (index, comment).

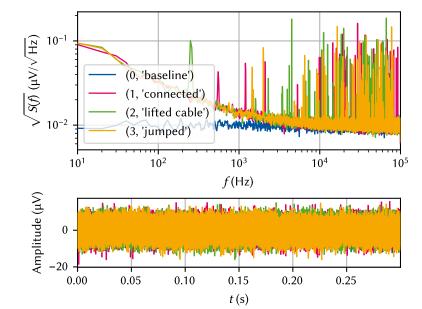


Figure 3.3: The python_spectrometer plot shown in Figure 3.2 when setting spect.plot_timetrace = True. This adds a subplot that shows the time series data from which the PSD was computed akin to what an oscilloscope would show. Note that this is the entire time series, *i.e.*, the data of length *L*, which is (by default, using Welch's method) segmented for spectrum estimation.

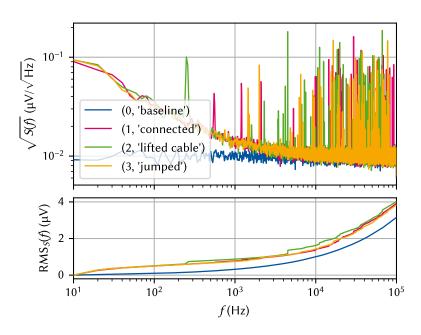


Figure 3.4: The python_spectrometer plot shown in Figure 3.2 when setting spect.plot_cumulative = True. This adds a subplot that shows the RMS (c.f. Equation 2.13) which can be helpful in evaluating the contribution of individual peaks in the spectrum to the total noise power. Both the oscilloscope subplot (Figure 3.3) and the RMS subplot can also be shown at the same time.

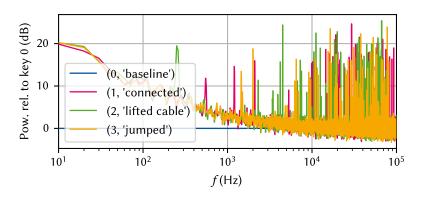


Figure 3.5: The python_spectrometer plot in relative mode. Starting from the state in Figure 3.2, we set spect.plot_j dB_scale = True as well as spect.j plot_amplitude = False and spect.j plot_density = False to compare the relative noise powers with respect to the baseline.

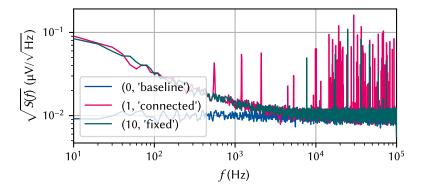


Figure 3.6: The python_spectrometer plot after multiple additional spectra were acquired and hidden. Hiding spectra that one is not interested in anymore is achieved through spect.jhide(*range(2, 10)). This is reversed by spect.show(*range(2, 10)). Data can also be dropped (spect.drop(key)) or deleted (spect.delete(key)) from the internal cache and disk, respectively.

The attribute plot_density controls whether the *power spectral density* or the *power spectrum*.¹⁷ Scaling the data to the power spectrum instead of the density, Bob can get an estimate of the RMS at a single frequency by reading off the peak height. Additionally displaying the data in dB then gives insight into relative noise powers of different spectra.

Bob carries on with his enterprise and continues to acquire spectra until, finally, he finds the source of his noise! Alas, his spectrometer plot is now overflowing with plotted data when really he just wants to compare the baseline, the original, noisy state, and the final, clean spectrum. He simply calls

```
spect.hide(*range(2, 10))
```

to hide the eight spectra of unsuccessful debugging, leaving him with a plot as shown in Figure 3.6.

Finally, happy with the results, Bob serializes the state of the spectrometer to disk, allowing him to pick up where he left off at a later point in time:

```
| spect.serialize_to_disk('2032-12-24_noise_hunting')
```

The next week, Bob is asked by his team about his progress on debugging the noise in their setup. Even though he is working from home that day and does not have access to the lab computer, Bob simply uses his laptop computer and pulls up the Spectrometer session stored on the server, allowing them to interactively discuss the spectra:

```
file = '2032-12-24_noise_hunting'
# Read-only instance because no DAQ attached
spect = Spectrometer.recall_from_disk(savepath / file)
```

This opens up the plot shown in Figure 3.6 again. While they cannot acquire new spectra in this state, ¹⁸ they can still use all the plotting features like showing or hiding spectra, or changing plot types as discussed above.

17: At this point, we *do* need to distinguish between the PSD and the power spectrum counter to sidenote 2 in Chapter 2. The PSD and power spectrum are related by the equivalent noise bandwidth (ENBW),

Spectral density
$$\xrightarrow{\times \text{ENBW}}$$
 Spectrum,

which is itself a function of the sampling rate and the properties of the spectral window [9],

ENBW =
$$f_s \frac{\sum_n \hat{w}_n^2}{\left[\sum_n \hat{w}_n\right]^2}$$
. (3.3)

18: They could of course always attach a DAQ instance to the spectrometer and continue as they were.

3.2.2 Live spectrum acquisition

Manually recording spectra in the workflow outlined in Subsection 3.2.1 becomes tedious at some point, and experimenters tend to become negligent with keeping metadata and comments up to date as they continue to change settings. Once a certain number of spectra has been obtained, the spectrometer plot also becomes crowded, and hiding old spectra manually is cumbersome. Moreover, each time a spectrum is captured, data is saved to disk, potentially accruing large amounts of disk space. For these

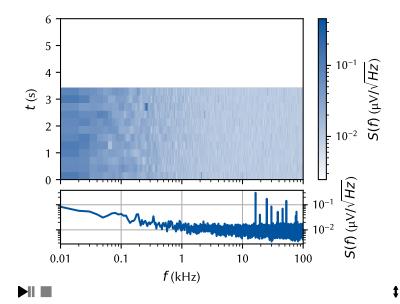


Figure 3.7: Spectrometer live view. The bottom plot shows the most recently acquired spectrum, while the top plot shows a waterfall plot of the most recent ones. Data acquisition runs in a background thread, keeping the interpreter responsive and available to interact with instruments, for example. The icons in the bottom left and right corner allow interacting with the live view.

reasons, the python_spectrometer package also offers a non-persistent live mode for displaying spectra continuously.

This mode is facilitated by the qutil.plotting.live_view module that provides asynchronous plotting functionality based on matplotlib. The live_view module supports both the multithreading and multiprocessing paradigms for concurrency in order to keep the interpreter responsive. In the former case, the window hosting the figure runs in the main thread and is kept responsive using matplotlib's GUI event loop mechanisms. In the latter, the plotting takes place in a separate process, resulting in true parallelism.

The live mode is started with the Spectrometer.live_view() method. Data is continuously acquired¹⁹ in a background thread using the same DAQ interface as the serial mode. Instead of saving the data on disk and managing plotting from python_spectrometer, however, the data is fed into a queue.²⁰ A live_view.IncrementalLiveView2D object then retrieves the data from the queue and handles plotting in a GUI event loop. To start a spectroscopy session with the same parameters as in Subsection 3.2.1 with a given Spectrometer object (c.f. Listing 3.7), we would call

view = spect.live_view(f_min=1e1, f_max=1e5, in_process=True)

which would open a figure window such as that shown in Figure 3.7. Similar to take(), the data acquisition parameters are passed to live_| view() as keyword arguments. The in_process argument specifies if multiprocessing (True) or multithreading (False) is used. Dictionaries with customization parameters for the live_view object can further be passed to the method.

19: Technically, a very large n_avg is used.

20: A queue is a concurrency mechanism for exchanging data between multiple threads or processes.

21: The difference is, of course, that we do not need to specify a comment since no data is retained.

Conclusion and outlook



N this part, I presented the python_spectrometer Python package for interactive, backend-agnostic spectrum acquisition. I first introduced the theory behind spectral noise estimation in Chapter 2. There, I discussed

Outlook. There are several possible avenues for future development of the python_spectrometer package. An obvious case is adding support for more DAQ hardware instruments by implementing DAQ interfaces. Modular devices such as those offered by QBLOX¹ and Quantum Machines² are on track to become the new standard in quantum technology labs. Implementing drivers for these instruments would benefit the adoption of both the instruments and the python_spectrometer package.

Next, incorporating noise spectroscopy into the standard measurement workflow of quantum device experiments would allow experimentalists to quickly gauge noise levels as they are performing measurements. If for some reason the noise changed³ the experimentalist could quickly obtain insight into the noise by analyzing the spectrum. In a client-server architecture, which is inherently asynchronous, such as Zurich Instrument's LabOne⁴ software, this is already possible using the web interface. But of course the strength of the python_spectrometer package stems from its capacity to be utilized in conjunction with any hardware instrument.

One way to implement such functionality would be to introduce a proxy DAQ subclass to be used together with the live mode presented in Subsection 3.2.2. This proxy class would serve as an interface to external measurement software and expose two attributes; first, a data queue, into which the external code could place arbitrary time series data that was obtained during some measurement, and second, a shared dictionary to hold acquisition parameters as these might change between measurements. Because the live view mode runs in the background, the external measurement framework could push data to the queue whenever new data was taken without obstructing the measurement workflow.

Listing 4.1 shows a template design for such a TeeDAQ class. The setup() method ignores the input parameters and instead obtains the current settings from the shared settings proxy. Similarly, instead of fetching data from an instrument itself, the acquire() method attempts to fetch data from the shared data_queue and blocks the thread if no data is present, thereby efficiently idling and consuming no resources unless triggered by the external caller. A measurement framework would then interact with the TeeDAQ object as exemplarized by the following code:

```
daq = TeeDAQ(...)
spect = Spectrometer(daq)
view = spect.live_view()
...
data = measure(fs, n_pts)
daq.settings.update(fs=fs, n_pts=n_pts)
daq.data_queue.put(data)
```

Measurement frameworks integrating with this interface could thus provide experimentalists live feedback on current noise levels with negligible overhead and minimal code adaptation.

```
1: https://www.qblox.com/research
2: https://www.quantum-machines.co/
```

- 3: As it happens often, unfortunately.
- 4: https://www.zhinst.com/ch/en/i nstruments/labone/labone-instrum ent-control-software

```
# dag/tee.py
import dataclasses
import threading
import multiprocessing as mp
from .base import DAQ
@dataclasses.dataclass
class TeeDAQ(DAQ):
    settings:

→ mp.managers.DictProxy

    data_queue: mp.JoinableQueue
    stop_event: threading.Event
    def setup(self, **_):
        settings = self.
        → DAQSettings(self.

    settings)

        return settings.

    to_consistent_dict()

    def acquire(self, **_):
        while not self.

    stop_event.is_set():

           yield self.

    get(block=True)
```

Listing 4.1: Template design for a proxy DAQ implementation to stream noise spectra from an external measurement framework. The settings attribute is a dictionary proxy shared between processes and used to pass acquisition parameters from the measurement framework to python_spectrometer.

- 5: Again, we use the two terms interchangably unless otherwise indicated, see sidenote 2 in Chapter 2.
- 6: Again assuming wide-sense stationary processes.

- 7: If one sticks to single DAQ instances managing single instruments.
- 8: Except possibly ensuring thread safety and timing synchronization if multiple DAQ objects communicate with the same physical instrument.

Listing 4.2: Proposed DAQ workflow for estimating cross-spectra. Each hardware channel (same or different instruments) is assigned to a DAQ object. After instrument configuration, it is asserted that the parameters match. Finally, data is fetched from both channels and fed into a CSD estimator. Note that triggering would need to be implemented externally.

[19]: (n.d.), Csd — SciPy v1.15.2 Manual
9: In practice, working with the normalized CSD, or correlation coefficient [20, 21]

$$r_{ij}(\omega) = \frac{S_{ij}(\omega)}{\sqrt{S_i(\omega)S_j(\omega)}}$$
 (4.2)

with $S_i(\omega)$ the PSD of process x_i would likely be more favorable.

10: It could be worthwile to add a subplot to display both the magnitude and phase of the complex quantity $S_{ij}(\omega)$.

improve this part.

Finally, it might be useful to not only allow estimating PSDs but also cross power spectral densitys (CSDs) or *cross-spectra*. The cross-spectrum⁵ is the Fourier transform of not the auto-correlation but the cross-correlation function $C(\tau)$ (*c.f.* Equation 2.3) between two random processes. Take a set of processes $\{x_1(t), x_2(t), \dots, x_n(t)\}$ that correspond to noise measured at different locations in a sample. The cross-correlation function between variables x_i and x_j is then given by⁶

$$C_{ij}(\tau) = \langle x_i(t)^* x_i(t+\tau) \rangle. \tag{4.1}$$

This function (and its Fourier pair the cross-spectrum $S_{ij}(\omega)$) quantifies the degree of correlation between noise at site i and noise at site j. Unlike the *auto*-spectrum (or self-spectrum), the cross-spectrum is always a complex quantity, even for real $x_i(t)$. It is not hard to see that for quantum processors, for example, these kinds of correlations could have significant impact on operation, and on error correction in particular [16–18]. To incorporate cross-spectra in the python_spectrometer package, only small changes should be necessary.

First, the data acquisition logic would need to be adapted. Two possible routes suggest themselves here; first, specialized DAQ classes could be implemented that, in place of yielding one batch of time series data, yield two batches each time they are queried. This approach first of all requires instruments with multiple channels, which is not necessarily given. Furthermore, it would incur additional coding efforts by having to reimplement each DAQ class for cross-spectra. On the other hand, it would arguably make synchronization between channels easier to achieve.

A less involved path would adapt the Spectrometer class to work with multiple DAQs. This would not involve additional driver work⁸ and allow the Spectrometer object to ensure synchronization between the different DAQs. The internal workflow shown in Listing 3.3 would then need to be slightly adapted to the code shown in Listing 4.2. The downside of this approach is that synchronization of different instruments or channels would need to be taken care of externally.

Further code adaptations would involve minor changes such as replacing the spectrum estimator with <code>scipy.signal.csd()</code> [19] for CSD estimation and make the plotting conform to complex data. 10

Part II

CHARACTERIZATION AND IMPROVEMENTS OF A MILLIKELVIN CONFOCAL MICROSCOPE

Part III

ELECTROSTATIC TRAPPING OF EXCITONS IN SEMICONDUCTOR MEMBRANES

Part IV

A FILTER-FUNCTION FORMALISM FOR UNITAL QUANTUM OPERATIONS



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Special Terms

```
A
API Application Programming Interface. 14
ASD amplitude spectral density. 15
C
CSD cross power spectral density. 19, 20
D
DAQ data acquisition device. 9, 11, 12, 14, 19
E
ENBW equivalent noise bandwidth. 17
F
FFT fast Fourier transform. 9
P
PSD power spectral density. 5–7, 9, 12, 13, 15–17, 19, 20
R
RMS root mean square. 6, 15–17
T
TIA transimpedance amplifier. 13
TLF two-level fluctuator. 5
```